APPENDIX A: SCOPE OF WORK

WOODARD & CURRAN
Engineering ▪ Science ▪ Operations

CORPORATE OFFICES: ME, MA, NH, CT, NY, GA, FL
Operational offices throughout the U.S.

MEMORANDUM

TO: Edward Logue, MEDEP

FROM: Karl Kasper, W&C

CC Mike Curtis, GP

DATE: May 30, 2006 (Version 2.0)

RE: SOW Updated 5/30/06 - Environmental Site Investigation of Georgia Pacific Paper Mill, Old Town, Maine

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The following is an updated Scope of Work (SOW), based on direction from the Maine Department of Environmental Protection (MEDEP) and results obtained during field work conducted 5/23 through 5/26. This memorandum describes the tasks completed and to be performed in conjunction with a Phase I and Phase II Environmental Site Assessment (ESA) investigation taking place at the Georgia Pacific (GP) paper mill. It is currently anticipated that draft Phase I and draft Phase II reports will be provided to the MEDEP by Friday June 2, 2006, however, as new data becomes available, updates will be provided in a timely manner.

Phase I

The Phase I ESA is being conducted using accepted industry standards in accordance with ASTM Standard E 1527-05 and 40 CFR Part 312, Standards and Practices for All Appropriate Inquiries (AAI) (November 1, 2005). The Phase I ESA is being conducted for the purpose of identifying information and/or the presence of a release, or disposal or threat of release, of potential hazardous materials affecting the Subject Property. The Phase I report will be formatted to include all required sections, as outlined in ASTM Standard E 1527-05.

Phase II

Based on the preliminary findings of the Phase I and direction provided by the MEDEP, the following SOW was proposed for the Phase II ESA. This scope will most likely be modified based on the findings of the Phase I or Phase II, at the direction of the MEDEP, or because of limitation imposed by surface and subsurface conditions. A supplemental investigation is being conducted at the south end of the GP property. This additional work is being conducted at the direction of the MEDEP. It is described in more detail below. The following is a brief summary of the Phase II investigation.

Task 1 - Utility Clearance

W&C, working with Maine DigSafe, will obtain a DigSafe permit number and clearance. It is our understanding that all on-site utilities are privately owned by GP. W&C will obtain direction from GP regarding on-site clearance for installation of borings and monitoring wells on the GP property. W&C

will direct and coordination with the drilling subcontractor. It will be the responsibility of GP to identify all on-site utilities.

Task 2 - Health and Safety Plan

A site-specific health and safety plan that provides emergency contact information and a planned route to the nearest health care facility will be provided by W&C. The health and safety plan will identify potential hazards at the job site and what measures will be taken to ensure protection from these hazards.

Task 3 - PCB Sampling

Based on historic records, 33 polychlorinated biphenyls (PCB) transformers were once present at the facility. According to GP, the facility is currently PCB free (e.g., all know PCB transformer have been changed out or removed). W&C will collect up to 35 soil, chip, or wipe samples from below or around these former transformer locations. Samples will be analyzed for PCBs at an off-site laboratory. Preliminary results should be available within 48-hours of sample submittal.

Task 4 - Soil Borings

Up to 50 GeoProbe soil borings are planned for installation around the mill property. These borings will be completed by two GeoProbe rigs. The locations of these soil borings will be determined in the field. Installation of the borings will be based on accessibility and site conditions. Based on conversations with the drilling contractor, the water table is expected at five to 10 feet below grade. The borings will extend approximately five feet below the water table, if possible. Subsurface soils will be screened visually or by a photoionization detector (PID) to identify potential contamination (see soil sampling below). In addition, a groundwater sample will be collected for on-site volatile organic compound (VOC) analysis from a GeoProbe soil boring. Groundwater sampling may be limited by subsurface conditions (e.g., bedrock).

All borings will be decommissioned, sealed with Bentonite pellets, and covered at the end of the investigation.

Task 4a - Subsurface GeoProbe Soil Sampling

Up to 15 soil samples will be collected from the soil borings. The location from which each of the soil borings will be collected will be based on PID readings or visual indications of contamination. The depth at which the highest PID reading is obtained or other visual indication of contamination will be the location from which the sample will be collected. Soil samples will be analyzed for VOCs, semi-volatile organic compounds (SVOCs), and RCRA 8 metals at the off-site laboratory, with preliminary results within 48-hours.

Task 4b - Subsurface GeoProbe Groundwater Sampling/On-Site Laboratory

At each of the soil borings that reach the water table, an attempt will be made to collect a groundwater sample for VOC analysis. It is anticipated that up to 50 groundwater samples will be analyzed by the on-site laboratory for VOC analysis.

A mobile laboratory will be setup at the site for VOC analysis. This will allow for 'real-time' analytical results to be used to direct subsequent sampling locations and depths. The on-site laboratory uses

standard laboratory equipment and methods. VOCs will be analyzed by a modified USEPA Method 8260.

In addition, up to five locations will also be sampled for SVOCs and RCRA-8 metal analysis at the off-site laboratory with preliminary result within 48-hour results.

Task 5 - Monitoring Well Installation

Based on the findings and the ability or inability of the GeoProbe rigs to obtain groundwater samples, sampled will be collected using a standard hollow stem auger drill rig to install up to four monitoring wells at the site. These wells will be constructed of 2-inch SCH 40 PVC screens and risers. Protective casing will be installed to protect these wells.

Task 5a - Groundwater Samples from Monitoring Wells

Groundwater samples will be collected from each of the monitoring wells these samples will be analyzed for VOAs, SVOCs and RCRA 8 metals analysis at the off-site laboratory with preliminary result within 48-hour results.

Task 6 - Analytical QA/QC

The following describes the quality assurance and quality control measures to be used to evaluate sample collection techniques and analytical methods.

- Equipment Blanks – an equipment blank will be collected for PCB sampling equipment.
- On-Site Laboratory – all samples analyzed at the on-site laboratory will be analyzed at a EPA Region 1 Tier I Level. This will include a duplicate sample at the rate of 1 in 10. Additional description of the on-site laboratory QA/QC program will be provided as an appendix to the Phase II report.
- Off-Site Laboratory – all samples analyzed at the on-site laboratory will be analyzed at a EPA Region 1 Tier II Level. This will include a duplicate sample at the rate of 1 in 10. Additional description of the on-site laboratory QA/QC program will be provided as an appendix to the Phase II report.
- Data Validation - because of time constraints to complete a Tier I or Tier II, data validation will not be completed prior to submitting the draft Phase II report. QC information will be provided from the laboratories such that these evaluations will be completed, as time allows

Task 7 - Phase II Summary Report

The findings of the Phase II investigation will be summarized in a letter report that includes sample results, field observations, and recommendations.

Task 8 - Schedule – It is currently anticipated that this SOW will be completed to the following schedule:

Updated Phase I Schedule

- DEP file review – May 19 and 22
- Preliminary Site Visit – May 19

- Second Site Visit – May 30
- Draft Report – June 2
- Final Report – June 9

Updated Phase II Schedule

- Utility Clearance – May 23 through 26
- PCB Sampling – May 23 and 24
- On-site laboratory and GeoProbe Investigation – May 24 through 26
- Drilling and groundwater sampling – May 25 and 26
- Off-site sample analyses – May 24 through June 1
- Draft Report – June 2
- Final Report – June 9

Task 9 – Phase II Amendment #1 – Based on information collected during the Phase I and Phase II investigation, MEDEP has directed W&C to conduct a supplemental investigation of Lot #2 at the south end of the GP property. This lot is approximately 67 acres and consists primarily of two former bark storage areas, a former sludge storage area, former lay-down area, and former gravel pit. The following is a summary of the proposed scope of work for this supplemental investigation.

- Former Northern Bark Storage Area – Based on conversations with current GP employees, this area was used to store bark. Based on historic topographic maps, this pile reached 30 to 40 feet above the ground surface. Currently, the majority of this material has been removed and was reportedly burned in the recovery boiler or sold as mulch. Current elevations in this area are approximately at the former ground surface. Because of a nesting bald eagle directly adjacent to this area, no subsurface explorations are planned. A 12/6/1990 spill report identifies this area as the site of a former PCB spill. W&C will collect two composite surface soil samples from the reported spill area. Each composite will be made up of four discrete samples. These samples will be submitted to the off-site laboratory for PBC analyses. Results should be available within 24-hours of submittal.

- Former Southern Storage Area – Based on conversations with current GP employees, this area served as bark and sludge storage area. The former southern storage area was located on Lot #2 to a height of 30 to 40 feet above the ground surface. To evaluate the nature and distribution of the remaining material, W&C is conducting a limited one day test pit investigation. This will include several test pits located around the edge of the former storage area to identify the edge of the former storage area.. Six subsurface soil samples will be collected from selected test pits for VOC, SVOC, and metals. Results should be available within 24-hours of submittal.

- Former Lay-down Area – A small area south west of the former southern storage area is a reported lay-down area for recycled equipment. W&C will conduct a limited test pitting in this area to evaluate if fill exist in this area, the extent and nature of fill in this area, if present. . Additional soil samples may be collected depending on visual observations.

- Former Gravel Pit – At the southern most end of Lot #2 is a former gavel pit, apparently used during past construction phases at the mill. While it is clear that the current landforms in this area are not natural, no test pitting or sampling is planned at this time.

APPENDIX B: BORING & TEST PIT LOGS

GeoProbe Logs
Monitoring Well Logs
Test Pit Logs

Appendix B
Geoprobe Logs
Georgia-Pacific Paper Mill
Old Town, Maine

WOODARD & CURRAN Engineering • Science • Operations					Geoprobe Log				Boring No.
					Georgia Pacific				SB-01
41 Hutchins Drive					Old Town, Maine				
Portland, ME 04102									
START DATE: May 25, 2006					BORING DEPTH: 10'				
CONTRACTOR: EPI					TEMPORARY WELL DEPTH: 10'				
DRILLER: Jason					TEMPORARY RISER LENGTH: N/A				
GEOLOGIST: Josh Bowe					TEMPORARY SCREEN LENGTH: N/A				
DRILLING METHOD: Direct Push					SLOT SIZE: 10				
Depth (feet)	**Sample Range**	**Sample Number**	**Rec. Length**	**PID (ppm)**	**Field Description and Remarks**				
0	0-4		2'		0-4'	f/m brown sands, some silt			
						Dark staining, slight odor			
5						Refusal at 10'			
10									
15									
20									
Groundwater Level Data									
Date	Time	Groundwater Depth (ft)			Summary				
						Soil Samples		1	
						Groundwater Samples		1	
						Soil Gas Samples		0	

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WOODARD & CURRAN
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41 Hutchins Drive
Portland, ME 04102

Geoprobe Log — **Boring No.**
Georgia Pacific — **SB-02**
Old Town, Maine

START DATE: May 24, 2006
CONTRACTOR: EPI
DRILLER: Jason
GEOLOGIST: Josh Bowe
DRILLING METHOD: Direct Push
BORING DEPTH: 4'
TEMPORARY WELL DEPTH: N/A
TEMPORARY RISER LENGTH: N/A
TEMPORARY SCREEN LENGTH: N/A
SLOT SIZE: N/A

Depth (feet)	Sample Range	Sample Number	Rec. Length	PID (ppm)	Field Description and Remarks
0	0-4		2'		0-4' f/m brown sands, some silt. Some fine gravel
					End of soil exploration at 4'
5					
10					
15					
20					

Groundwater Level Data

Date	Time	Groundwater Depth (ft)

Summary

Soil Samples	1
Groundwater Samples	0
Soil Gas Samples	0

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41 Hutchins Drive
Portland, ME 04102

Geoprobe Log — **Boring No.**
Georgia Pacific — **SB-03**
Old Town, Maine

START DATE: May 24, 2006
CONTRACTOR: EPI
DRILLER: Jason
GEOLOGIST: Josh Bowe
DRILLING METHOD: Direct Push
BORING DEPTH: 4'
TEMPORARY WELL DEPTH: N/A
TEMPORARY RISER LENGTH: N/A
TEMPORARY SCREEN LENGTH: N/A
SLOT SIZE: N/A

Depth (feet)	Sample Range	Sample Number	Rec. Length	PID (ppm)	Field Description and Remarks
0	0-4		1.5'		0-4' f/m brown sands, some silt. Some fine gravel
					End of soil exploration at 4'
5					
10					
15					
20					

Groundwater Level Data

Date	Time	Groundwater Depth (ft)

Summary

Soil Samples	1
Groundwater Samples	0
Soil Gas Samples	0

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Georgia-Pacific Paper Mill
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					Geoprobe Log				**Boring No.**
					Georgia Pacific				**SB-04**
41 Hutchins Drive					**Old Town, Maine**				
Portland, ME 04102									
START DATE: May 24, 2006					BORING DEPTH: 4'				
CONTRACTOR: EPI					TEMPORARY WELL DEPTH: N/A				
DRILLER: Jason					TEMPORARY RISER LENGTH: N/A				
GEOLOGIST: Josh Bowe					TEMPORARY SCREEN LENGTH: N/A				
DRILLING METHOD: Direct Push					SLOT SIZE: N/A				
Depth	**Sample**	**Sample**	**Rec.**	**PID**					
(feet)	**Range**	**Number**	**Length**	**(ppm)**	**Field Description and Remarks**				
0	0-4		1'		0-4'	f/m brown sands, some silt. Some fine gravel			
						End of soil exploration at 4'			
5									
10									
15									
20									
Groundwater Level Data									
Date	Time	Groundwater Depth (ft)			Summary				
						Soil Samples		1	
						Groundwater Samples		0	
						Soil Gas Samples		0	

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Georgia-Pacific Paper Mill
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WOODARD & CURRAN Engineering • Science • Operations					Geoprobe Log				Boring No.
					Georgia Pacific				SB-05
41 Hutchins Drive					Old Town, Maine				
Portland, ME 04102									
START DATE: May 25, 2006					BORING DEPTH: 6'				
CONTRACTOR: EPI					TEMPORARY WELL DEPTH: 6'				
DRILLER: Jason					TEMPORARY RISER LENGTH: N/A				
GEOLOGIST: Josh Bowe					TEMPORARY SCREEN LENGTH: N/A				
DRILLING METHOD: Direct Push					SLOT SIZE: N/A				
Depth	**Sample**	**Sample**	**Rec.**	**PID**					
(feet)	**Range**	**Number**	**Length**	**(ppm)**	**Field Description and Remarks**				
0	0-4		2.5'		0-4'	f/m brown sands, fine gravel, some silt.			
	4-6		6"		4-6'	f/m brown sands, fine gravel, some silt.			
5									
						Refusal at 6' bgs			
10									
15									
20									
Groundwater Level Data									
Date	Time	Groundwater Depth (ft)			Summary				
						Soil Samples		2	full suite (0-4')
						Groundwater Samples		No Water for Sampling	
						Soil Gas Samples		0	

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Georgia-Pacific Paper Mill
Old Town, Maine

WOODARD & CURRAN Engineering • Science • Operations					Geoprobe Log				Boring No.
					Georgia Pacific				SB-06
41 Hutchins Drive					Old Town, Maine				
Portland, ME 04102									
START DATE: May 25, 2006					BORING DEPTH: 2'				
CONTRACTOR: EPI					TEMPORARY WELL DEPTH: N/A				
DRILLER: Jason					TEMPORARY RISER LENGTH: N/A				
GEOLOGIST: Josh Bowe					TEMPORARY SCREEN LENGTH: N/A				
DRILLING METHOD: Direct Push					SLOT SIZE: N/A				
Depth	**Sample**	**Sample**	**Rec.**	**PID**					
(feet)	**Range**	**Number**	**Length**	**(ppm)**	**Field Description and Remarks**				
0	0-2		6"		0-2'	f/m brown sands, f/m gravel, trace silt. Mostly fill.			
						Refusal at 2' bgs			
5									
10									
15									
20									
Groundwater Level Data									
Date	Time	Groundwater Depth (ft)			Summary				
						Soil Samples		1	
						Groundwater Samples		No Water for Sampling	
						Soil Gas Samples		0	

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WOODARD & CURRAN Engineering • Science • Operations					Geoprobe Log		Boring No.
					Georgia - Pacific		SB-10
41 Hutchins Drive					Old Town, Maine		
Portland, ME 04102							
START DATE: May 24, 2006					BORING DEPTH: 1.33 feet		
CONTRACTOR: Environmental Services Network					TEMPORARY WELL DEPTH: N/A		
DRILLER: David Dionne					TEMPORARY RISER LENGTH: N/A		
GEOLOGIST: Danielle Taylor					TEMPORARY SCREEN LENGTH: N/A		
DRILLING METHOD: Direct Push					SLOT SIZE: N/A		
Depth (feet)	**Sample Range**	**Sample Number**	**Rec. Length**	**PID (ppm)**	**Field Description and Remarks**		
0	0-1.33		0.67		0 to 0.5	Reddish brown fine to medium SAND	
					0.5 to 0.67	Reddish brown fine to medium SAND with some black material, slight odor.	
5						Refusal at 1.33' bgs	
10							
15							
20							
Groundwater Level Data							
Date	Time	Groundwater Depth (ft)			Summary		
					Soil Samples	1	
					Groundwater Samples	0	
					Soil Gas Samples	0	

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Old Town, Maine

WOODARD & CURRAN Engineering • Science • Operations					Geoprobe Log		Boring No.
					Georgia - Pacific		SB-11
41 Hutchins Drive					Old Town, Maine		
Portland, ME 04102							
START DATE: May 24, 2006					BORING DEPTH: 12'		
CONTRACTOR: Environmental Services Network					TEMPORARY WELL DEPTH: 10'		
DRILLER: David Dionne					TEMPORARY RISER LENGTH: N/A		
GEOLOGIST: Danielle Taylor					TEMPORARY SCREEN LENGTH: 4'		
DRILLING METHOD: Direct Push					SLOT SIZE: N/A		
Depth (feet)	**Sample Range**	**Sample Number**	**Rec. Length**	**PID (ppm)**	**Field Description and Remarks**		
0	0-4		3		0 to 1	Brown fine to coarse SAND	
					1 to 2	Fine to medium dark brown to black SAND with some	
						black SILT and CLAY, strong odor.	
					2 to 3	Some wood chips to black fine to medium SAND,	
	4-8		4			strong odor.	
5					4 to 7	Strong to slight odor, wood chips, fine to medium	
						black/dk gray SAND	
					7 to 7.5	Reddish brown fine SAND, some medium SAND,	
	8-12		4			some silt, wet.	
					7.5 to 8	Reddish brown coarse SAND, some silt and fine	
10						brown SAND, more water.	
					8 to 8.5	Brown coarse to medium SAND	
					8.5-12	Brown clay, traces of orange medium SAND	
						End of exploration at 12 ft bgs	
15							
20							
Groundwater Level Data							
Date	Time	Groundwater Depth (ft)			Summary		
						Soil Samples	1
						Groundwater Samples	1
						Soil Gas Samples	0

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						Geoprobe Log				**Boring No.**
						Georgia - Pacific				**SB-12**
41 Hutchins Drive						**Old Town, Maine**				
Portland, ME 04102										
START DATE: May 24, 2006						BORING DEPTH: 7.5'				
CONTRACTOR: Environmental Services Network						TEMPORARY WELL DEPTH: 7.5'				
DRILLER: David Dionne						TEMPORARY RISER LENGTH: N/A				
GEOLOGIST: Danielle Taylor						TEMPORARY SCREEN LENGTH: 4'				
DRILLING METHOD: Direct Push						SLOT SIZE: N/A				
Depth (feet)	**Sample Range**	**Sample Number**	**Rec. Length**	**PID (ppm)**		**Field Description and Remarks**				
0	0-4		4		0-0.5	Asphalt				
					0.5 to 2	Brown medium to coarse SAND				
					2 to 3	Brown to reddish brown coarse SAND				
					3 to 4	Brown fine to coarse SAND, reddish SAND @ 3.75'				
	4-8		2			bgs, slight odor.				
5					4 to 7.5	Brown to reddish brown medium to coarse SAND				
						Refusal at 7.5' bgs				
10										
15										
20										
Groundwater Level Data										
Date	Time	Groundwater Depth (ft)				Summary				
						Soil Samples		1		
						Groundwater Samples		1		
						Soil Gas Samples		0		

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WOODARD & CURRAN Engineering • Science • Operations					Geoprobe Log		Boring No.
					Georgia - Pacific		SB-13
41 Hutchins Drive					Old Town, Maine		
Portland, ME 04102							
START DATE: May 24, 2006					BORING DEPTH: 10'		
CONTRACTOR: Environmental Services Network					TEMPORARY WELL DEPTH: 10'		
DRILLER: David Dionne					TEMPORARY RISER LENGTH: N/A		
GEOLOGIST: Danielle Taylor					TEMPORARY SCREEN LENGTH: 4'		
DRILLING METHOD: Direct Push					SLOT SIZE: N/A		
Depth (feet)	**Sample Range**	**Sample Number**	**Rec. Length**	**PID (ppm)**	**Field Description and Remarks**		
0	0-4		4		0-0.5	Asphalt	
					0.5 to 1	Reddish brown to brown coarse SAND	
					1 to 1.3	Black medium to coarse SAND, strong odor	
					1.3 to 2	Orange to reddish medium to fine SAND	
					2 to 2.5	Black fine to medium SAND	
5					2.5 to 3	Light brown medium to coarse SAND	
					3 to 3.5	Orange (0.2' thick) to Black medium to coarse SAND	
					3.5 to 4	Brown wood-like material	
						End of soil exploration at 4'	
10							
15							
20							
Groundwater Level Data							
Date	Time	Groundwater Depth (ft)			Summary		
						Soil Samples	1
						Groundwater Samples	1
						Soil Gas Samples	0

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					Geoprobe Log				**Boring No.**
					Georgia - Pacific				**SB-14**
41 Hutchins Drive					**Old Town, Maine**				
Portland, ME 04102									
START DATE: May 24, 2006					BORING DEPTH: 10'				
CONTRACTOR: Environmental Services Network					TEMPORARY WELL DEPTH: 10'				
DRILLER: David Dionne					TEMPORARY RISER LENGTH: N/A				
GEOLOGIST: Danielle Taylor					TEMPORARY SCREEN LENGTH: 4'				
DRILLING METHOD: Direct Push					SLOT SIZE: N/A				
Depth (feet)	**Sample Range**	**Sample Number**	**Rec. Length**	**PID (ppm)**	**Field Description and Remarks**				
0	0-4		2.5		0-0.5	Wood-like material			
					0.5 to 1	Brown medium to coarse SAND			
					1 to 1.5	Brown medium to coarse SAND, some staining			
						material			
					1.5 to 2.5	Brown fine to very fine SAND,			
5						brown fine SAND and silt in drive-shoe.			
						End of soil exploration at 4' bgs			
10									
15									
20									
Groundwater Level Data									
Date	Time	Groundwater Depth (ft)			Summary				
						Soil Samples		0	
						Groundwater Samples		1	
						Soil Gas Samples		0	

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						Geoprobe Log				**Boring No.**
						Georgia - Pacific				**SB-15**
41 Hutchins Drive						**Old Town, Maine**				
Portland, ME 04102										
START DATE: May 24, 2006						BORING DEPTH: 8'				
CONTRACTOR: Environmental Services Network						TEMPORARY WELL DEPTH: 8'				
DRILLER: David Dionne						TEMPORARY RISER LENGTH: N/A				
GEOLOGIST: Danielle Taylor						TEMPORARY SCREEN LENGTH: 5'				
DRILLING METHOD: Direct Push						SLOT SIZE: N/A				
Depth	**Sample**	**Sample**	**Rec.**	**PID**						
(feet)	**Range**	**Number**	**Length**	**(ppm)**	**Field Description and Remarks**					
0	0-4		4		0 to 1	Brown coarse SAND				
					1 to 2	Brown coarse SAND, some fine to very fine brown				
						SAND and silt				
					2 to 3	Brown to gray coarse SAND, some fine to very fine				
	4-8		4			gray SAND and silt				
5					3 to 4	Gray fine to very fine SAND and silt, trace of				
						coarse SAND, strong odor				
					4 to 5	Light gray fine to very fine SAND, strong odor				
					5 to 8	Black wood-like material, slight odor				
10						End of soil investigation at 8' bgs				
15										
20										
Groundwater Level Data										
Date	Time	Groundwater Depth (ft)			Summary					
						Soil Samples		2		
						Groundwater Samples		2		
						Soil Gas Samples		0		

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WOODARD & CURRAN Engineering • Science • Operations					Geoprobe Log				Boring No.
					Georgia - Pacific				SB-16
41 Hutchins Drive					Old Town, Maine				
Portland, ME 04102									
START DATE: May 24, 2006					BORING DEPTH: 1'				
CONTRACTOR: Environmental Services Network					TEMPORARY WELL DEPTH:				1'
DRILLER: David Dionne					TEMPORARY RISER LENGTH: N/A				
GEOLOGIST: Danielle Taylor					TEMPORARY SCREEN LENGTH: N/A				
DRILLING METHOD: Direct Push					SLOT SIZE: N/A				
Depth (feet)	**Sample Range**	**Sample Number**	**Rec. Length**	**PID (ppm)**	**Field Description and Remarks**				
0	0-1		1		0 to 0.5	Wood-like material			
					0.5 to 1	Brown coarse to medium SAND			
						Refusal at 1' bgs			
5									
10									
15									
20									
Groundwater Level Data									
Date	Time	Groundwater Depth (ft)			Summary				
						Soil Samples		0	
					Groundwater Samples			0	
					Soil Gas Samples			0	

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WOODARD & CURRAN Engineering • Science • Operations					Geoprobe Log		Boring No.
					Georgia - Pacific		SB-17
41 Hutchins Drive					Old Town, Maine		
Portland, ME 04102							
START DATE: May 24, 2006					BORING DEPTH: 7.5'		
CONTRACTOR: Environmental Services Network					TEMPORARY WELL DEPTH: 7.5'		
DRILLER: David Dionne					TEMPORARY RISER LENGTH: N/A		
GEOLOGIST: Danielle Taylor					TEMPORARY SCREEN LENGTH: 4'		
DRILLING METHOD: Direct Push					SLOT SIZE: N/A		
Depth (feet)	**Sample Range**	**Sample Number**	**Rec. Length**	**PID (ppm)**	**Field Description and Remarks**		
0	0-4		1.5		0 to 0.5	Brown coarse angular SAND	
					0.5 to 1.5	Brown medium SAND, some fine SAND and silt	
5						End of soil exploration at 4' bgs	
10							
15							
20							

Groundwater Level Data			Summary		
Date	Time	Groundwater Depth (ft)			
			Soil Samples	0	
			Groundwater Samples	1	
			Soil Gas Samples	0	

Appendix B
Geoprobe Logs
Georgia-Pacific Paper Mill
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WOODARD & CURRAN
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						Geoprobe Log				**Boring No.**
						Georgia - Pacific				**SB-18**
41 Hutchins Drive						**Old Town, Maine**				
Portland, ME 04102										
START DATE: May 24, 2006						BORING DEPTH: 14'				
CONTRACTOR: Environmental Services Network						TEMPORARY WELL DEPTH: 14'				
DRILLER: David Dionne						TEMPORARY RISER LENGTH: N/A				
GEOLOGIST: Danielle Taylor						TEMPORARY SCREEN LENGTH: 4'				
DRILLING METHOD: Direct Push						SLOT SIZE: N/A				
Depth (feet)	**Sample Range**	**Sample Number**	**Rec. Length**	**PID (ppm)**		**Field Description and Remarks**				
0	0-4		3'		0 to 0.5	Topsoil to light gray coarse SAND				
					0.5 to 1	Dark brown medium SAND, some black staining				
					1 to 2.5	Dark brown medium SAND, more black staining				
					2.5 to 3	Light brown fine sand and silt, some orange staining				
5						End of soil investigation at 4' bgs				
10										
15										
20										
Groundwater Level Data										
Date	Time	Groundwater Depth (ft)				Summary				
						Soil Samples		1		
						Groundwater Samples		1		
						Soil Gas Samples		0		

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Georgia-Pacific Paper Mill
Old Town, Maine

WOODARD & CURRAN Engineering • Science • Operations					Geoprobe Log		Boring No.
					Georgia - Pacific		SB-19
41 Hutchins Drive					Old Town, Maine		
Portland, ME 04102							
START DATE: May 24, 2006					BORING DEPTH: 10'		
CONTRACTOR: Environmental Services Network					TEMPORARY WELL DEPTH: 10'		
DRILLER: David Dionne					TEMPORARY RISER LENGTH: N/A		
GEOLOGIST: Danielle Taylor					TEMPORARY SCREEN LENGTH: 4'		
DRILLING METHOD: Direct Push					SLOT SIZE: N/A		
Depth (feet)	**Sample Range**	**Sample Number**	**Rec. Length**	**PID (ppm)**	**Field Description and Remarks**		
0	0-4		1'		0 to 0.5	Dark brown coarse SAND	
					0.5 to 1	Dark brown coarse SAND, some staining, slight odor	
						Rock in drive shoe	
5						End of soil investigation at 4' bgs	
10							
15							
20							
Groundwater Level Data							
Date	Time	Groundwater Depth (ft)			Summary		
					Soil Samples	0	
					Groundwater Samples	2	(2-6 bgs) and (6-10 bgs)
					Soil Gas Samples	0	

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Geoprobe Logs
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						Geoprobe Log				**Boring No.**
						Georgia - Pacific				**SB-20**
41 Hutchins Drive						**Old Town, Maine**				
Portland, ME 04102										
START DATE: May 24, 2006						BORING DEPTH: 14'				
CONTRACTOR: Environmental Services Network						TEMPORARY WELL DEPTH: 14'				
DRILLER: David Dionne						TEMPORARY RISER LENGTH: N/A				
GEOLOGIST: Danielle Taylor						TEMPORARY SCREEN LENGTH: 4'				
DRILLING METHOD: Direct Push						SLOT SIZE: N/A				
Depth (feet)	**Sample Range**	**Sample Number**	**Rec. Length**	**PID (ppm)**		**Field Description and Remarks**				
0	0-4		4'		0 to 0.5	Brown coarse SAND				
					0.5 to 1	Brown fine to very fine SAND and silt,				
					1 to 2	Light brown fine to very fine SAND and silt				
					2 to 4	Light brown fine to very fine SAND, more silt				
5						End of soil investigation at 4' bgs				
10										
15										
20										
Groundwater Level Data										
Date	Time	Groundwater Depth (ft)				Summary				
						Soil Samples		0		
						Groundwater Samples		1		
						Soil Gas Samples		0		

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					Geoprobe Log				**Boring No.**
					Georgia Pacific				**SB-21**
41 Hutchins Drive					**Old Town, Maine**				
Portland, ME 04102									
START DATE: May 25, 2006					BORING DEPTH: 14'				
CONTRACTOR: EPI					TEMPORARY WELL DEPTH: 14'				
DRILLER: Jason					TEMPORARY RISER LENGTH: N/A				
GEOLOGIST: Josh Bowe					TEMPORARY SCREEN LENGTH: N/A				
DRILLING METHOD: Direct Push					SLOT SIZE: 10				
Depth	**Sample**	**Sample**	**Rec.**	**PID**					
(feet)	**Range**	**Number**	**Length**	**(ppm)**	**Field Description and Remarks**				
0	0-4		1'		0-4'	f/m brown sands, f/m gravel, trace silt.			
	4-8		1'		4-8'	f/m brown sands, f/m gravel, trace silt.			
5									
10									
						Refusal at 14' bgs			
15									
20									
Groundwater Level Data									
Date	Time	Groundwater Depth (ft)			Summary				
						Soil Samples		2	full suite (0-4')
						Groundwater Samples		No Water for Sampling	
						Soil Gas Samples		0	

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WOODARD & CURRAN Engineering • Science • Operations					Geoprobe Log		Boring No.
					Georgia - Pacific		SB-22
41 Hutchins Drive					Old Town, Maine		
Portland, ME 04102							
START DATE: May 25, 2006					BORING DEPTH: 16'		
CONTRACTOR: Environmental Services Network					TEMPORARY WELL DEPTH: 16'		
DRILLER: David Dionne					TEMPORARY RISER LENGTH: N/A		
GEOLOGIST: Danielle Taylor					TEMPORARY SCREEN LENGTH: N/A		
DRILLING METHOD: Direct Push					SLOT SIZE: N/A		
Depth (feet)	**Sample Range**	**Sample Number**	**Rec. Length**	**PID (ppm)**	**Field Description and Remarks**		
0	0-4		3		0 to 1	Asphalt, Brown fine SAND, trace coarse SAND, dark brown silt	
					1 to 3	Brown fine SAND, dark brown silt	
	4-8		3		4 to 7	Brown fine SAND, some coarse to medium SAND	
5							
	8-12		4		8 to 10	Dark brown medium to coarse SAND, some silt	
10					10 to 12	Brown to gray coarse SAND, some GRAVEL	
	12-16		4		12 to 16	Gray to brown silt and clay, dry	
15							
20							
Groundwater Level Data							
Date	Time	Groundwater Depth (ft)			Summary		
					Soil Samples	1	
					Groundwater Samples	0	
					Soil Gas Samples	0	

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WOODARD & CURRAN Engineering • Science • Operations					Geoprobe Log				Boring No.
					Georgia Pacific				SB-23
41 Hutchins Drive					Old Town, Maine				
Portland, ME 04102									
START DATE: May 25, 2006					BORING DEPTH: 13'				
CONTRACTOR: EPI					TEMPORARY WELL DEPTH: 13'				
DRILLER: Jason					TEMPORARY RISER LENGTH: N/A				
GEOLOGIST: Josh Bowe					TEMPORARY SCREEN LENGTH: N/A				
DRILLING METHOD: Direct Push					SLOT SIZE: 10				
Depth (feet)	**Sample Range**	**Sample Number**	**Rec. Length**	**PID (ppm)**	**Field Description and Remarks**				
0	0-4		1'		0-4'	f/m brown sands and silt, f/m gravel. Mostly fill			
5									
10									
						Refusal at 13' bgs			
15									
20									
Groundwater Level Data					Summary				
Date	Time	Groundwater Depth (ft)			Soil Samples		1		
					Groundwater Samples		No Water for Sampling		
					Soil Gas Samples		0		

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					Geoprobe Log				**Boring No.**
					Georgia Pacific				**SB-24**
41 Hutchins Drive					**Old Town, Maine**				
Portland, ME 04102									
START DATE: May 25, 2006					BORING DEPTH: 8'				
CONTRACTOR: EPI					TEMPORARY WELL DEPTH: 8'				
DRILLER: Jason					TEMPORARY RISER LENGTH: N/A				
GEOLOGIST: Josh Bowe					TEMPORARY SCREEN LENGTH: N/A				
DRILLING METHOD: Direct Push					SLOT SIZE: 10				
Depth	**Sample**	**Sample**	**Rec.**	**PID**					
(feet)	**Range**	**Number**	**Length**	**(ppm)**	**Field Description and Remarks**				
0	0-4		2'		0-4'	f/m brown sands and silt, f/m gravel. Mostly fill			
						Some visual staining			
5									
						Refusal at 8' bgs			
10									
15									
20									
Groundwater Level Data									
Date	Time	Groundwater Depth (ft)			Summary				
						Soil Samples		1	
						Groundwater Samples		No Water for Sampling	
						Soil Gas Samples		0	

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						Geoprobe Log				**Boring No.**
						Georgia Pacific				**SB-25**
41 Hutchins Drive						**Old Town, Maine**				
Portland, ME 04102										
START DATE: May 25, 2006						BORING DEPTH: 11'				
CONTRACTOR: EPI						INSTALLED WELL DEPTH: 11'				
DRILLER: Jason						TEMPORARY RISER LENGTH: N/A				
GEOLOGIST: Josh Bowe						TEMPORARY SCREEN LENGTH: 4'				
DRILLING METHOD: Direct Push						SLOT SIZE: 10				
Depth	**Sample**	**Sample**	**Rec.**	**PID**						
(feet)	**Range**	**Number**	**Length**	**(ppm)**	**Field Description and Remarks**					
0	0-4		10"		0-4'	f/m brown sands, some brown silt.				
5										
10										
						Refusal at 11'				
15										
20										
Groundwater Level Data										
Date	Time	Groundwater Depth (ft)			Summary					
						Soil Samples		1		
						Groundwater Samples		4		(7-11') bgs
						Soil Gas Samples		0		

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					Geoprobe Log				**Boring No.**
					Georgia Pacific				**SB-26**
41 Hutchins Drive					**Old Town, Maine**				
Portland, ME 04102									
START DATE: May 25, 2006					BORING DEPTH: 12'				
CONTRACTOR: EPI					INSTALLED WELL DEPTH: 12'				
DRILLER: Jason					TEMPORARY RISER LENGTH: N/A				
GEOLOGIST: Josh Bowe					TEMPORARY SCREEN LENGTH: 4'				
DRILLING METHOD: Direct Push					SLOT SIZE: 10				
Depth (feet)	**Sample Range**	**Sample Number**	**Rec. Length**	**PID (ppm)**	**Field Description and Remarks**				
0	0-4		1'		0-4'	f/m brown sands, some brown silt.			
5									
						Groundwater Depth ~ 8' bgs			
10									
						Refusal at 12'			
15									
20									
Groundwater Level Data									
Date	Time	Groundwater Depth (ft)			Summary				
						Soil Samples		1	
						Groundwater Samples		2	
						Soil Gas Samples		0	

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						Geoprobe Log				**Boring No.**
						Georgia Pacific				**SB-27**
41 Hutchins Drive						**Old Town, Maine**				
Portland, ME 04102										
START DATE: May 25, 2006						BORING DEPTH: 8'				
CONTRACTOR: EPI						TEMPORARY WELL DEPTH:				8'
DRILLER: Jason						TEMPORARY RISER LENGTH: N/A				
GEOLOGIST: Josh Bowe						TEMPORARY SCREEN LENGTH: 4'				
DRILLING METHOD: Direct Push						SLOT SIZE: 10				
Depth (feet)	**Sample Range**	**Sample Number**	**Rec. Length**	**PID (ppm)**		**Field Description and Remarks**				
0	0-4		1'		0-4'	f/m brown sands, some brown silt.				
						slight staining and odor				
5										
						Refusal at 8'				
10										
15										
20										
Groundwater Level Data										
Date	Time	Groundwater Depth (ft)				Summary				
						Soil Samples		1		
						Groundwater Samples		No Water (4-8')		
						Soil Gas Samples		0		

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					Geoprobe Log				**Boring No.**
					Georgia Pacific				**SB-28**
41 Hutchins Drive					**Old Town, Maine**				
Portland, ME 04102									
START DATE: May 24, 2006					BORING DEPTH: 9'				
CONTRACTOR: EPI					TEMPORARY WELL DEPTH: 9'				
DRILLER: Jason					TEMPORARY RISER LENGTH: N/A				
GEOLOGIST: Josh Bowe					TEMPORARY SCREEN LENGTH: N/A				
DRILLING METHOD: Direct Push					SLOT SIZE: 10				
Depth (feet)	**Sample Range**	**Sample Number**	**Rec. Length**	**PID (ppm)**	**Field Description and Remarks**				
0	0-4'		8"		0-4'	f/m brown sands, some silt. Trace fine gravel.			
5									
						Refusal at 9' bgs			
10									
15									
20									
Groundwater Level Data									
Date	Time	Groundwater Depth (ft)			Summary				
						Soil Samples		1	
					Groundwater Samples		No Water (9-5')		
					Soil Gas Samples			0	

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41 Hutchins Drive
Portland, ME 04102

Geoprobe Log — **Boring No.**
Georgia Pacific — **SB-29**
Old Town, Maine

START DATE: May 24, 2006
CONTRACTOR: EPI
DRILLER: Jason
GEOLOGIST: Josh Bowe
DRILLING METHOD: Direct Push
BORING DEPTH: 4'
TEMPORARY WELL DEPTH: N/A
TEMPORARY RISER LENGTH: N/A
TEMPORARY SCREEN LENGTH: N/A
SLOT SIZE: N/A

Depth (feet)	Sample Range	Sample Number	Rec. Length	PID (ppm)	Field Description and Remarks
0	0-4		1'		0-4' f/m brown sands, some silt
					Slight odor and staining
					End of boring at 4'
5					
10					
15					
20					

Groundwater Level Data

Date	Time	Groundwater Depth (ft)

Summary

Soil Samples	1
Groundwater Samples	0
Soil Gas Samples	0

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WOODARD & CURRAN Engineering • Science • Operations					Geoprobe Log		Boring No.
					Georgia - Pacific		SB-30
41 Hutchins Drive					Old Town, Maine		
Portland, ME 04102							
START DATE: May 25, 2006					BORING DEPTH: 10'		
CONTRACTOR: Environmental Services Network					TEMPORARY WELL DEPTH: 10'		
DRILLER: David Dionne					TEMPORARY RISER LENGTH: N/A		
GEOLOGIST: Danielle Taylor					TEMPORARY SCREEN LENGTH: 4'		
DRILLING METHOD: Direct Push					SLOT SIZE: N/A		
Depth (feet)	**Sample Range**	**Sample Number**	**Rec. Length**	**PID (ppm)**	**Field Description and Remarks**		
0	0-4		3		0 to 0.5	Topsoil, gray puddy-like substance, brown coarse to fine SAND.	
					0.5 to1	Organics, dark brown to black medium to fine SAND trace gray coarse SAND	
					1 to 2	Dark brown medium to fine SAND, @ 2.5 change reddish brown coarse angular SAND	
5					2 to 3	Wood-like material, light brown fine to very fine SAND	
						End of soil investigation at 4' bgs	
10							
15							
20							
Groundwater Level Data							
Date	Time	Groundwater Depth (ft)			Summary		
						Soil Samples: 1	
						Groundwater Samples: 1	
						Soil Gas Samples: 0	

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					Geoprobe Log				**Boring No.**
					Georgia - Pacific				**SB-31**
41 Hutchins Drive					**Old Town, Maine**				
Portland, ME 04102									
START DATE: May 25, 2006					BORING DEPTH: 7.8'				
CONTRACTOR: Environmental Services Network					TEMPORARY WELL DEPTH: 7.8'				
DRILLER: David Dionne					TEMPORARY RISER LENGTH: N/A				
GEOLOGIST: Danielle Taylor					TEMPORARY SCREEN LENGTH: 4'				
DRILLING METHOD: Direct Push					SLOT SIZE: N/A				
Depth	**Sample**	**Sample**	**Rec.**	**PID**					
(feet)	**Range**	**Number**	**Length**	**(ppm)**	**Field Description and Remarks**				
0	0-4		2.5		0 to 1	Organics, brown fine to coarse SAND			
					1 to 2.5	Wood-like material, black material, slight odor			
	4-7.8'		4		4 to 7.8	Black coarse SAND, slight odor			
5						Weathered bedrock in drive shoe			
10						Refusal at 7.8' bgs			
15									
20									

Groundwater Level Data									
Date	Time	Groundwater Depth (ft)			Summary				
						Soil Samples		1	
					Groundwater Samples			1	
					Soil Gas Samples			0	

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						Geoprobe Log				**Boring No.**
						Georgia - Pacific				**SB-32**
41 Hutchins Drive						**Old Town, Maine**				
Portland, ME 04102										
START DATE: May 25, 2006						BORING DEPTH: 10'				
CONTRACTOR: Environmental Services Network						TEMPORARY WELL DEPTH: 10'				
DRILLER: David Dionne						TEMPORARY RISER LENGTH: N/A				
GEOLOGIST: Danielle Taylor						TEMPORARY SCREEN LENGTH: 4'				
DRILLING METHOD: Direct Push						SLOT SIZE: N/A				
Depth	**Sample**	**Sample**	**Rec.**	**PID**						
(feet)	**Range**	**Number**	**Length**	**(ppm)**	**Field Description and Remarks**					
0	0-4		3'		0 to 0.5	Asphalt				
					0.5 to 1	Brown medium SAND				
					1 to 2	Coarse angular GRAVEL				
					2 to 3	Brown medium to fine SAND, some coarse SAND				
5						End of soil investigation at 4' bgs				
					* sheen and strong odor in water sample (6-10' bgs)					
10										
15										
20										
Groundwater Level Data										
Date	Time	Groundwater Depth (ft)			Summary					
						Soil Samples		0		
						Groundwater Samples		1		
						Soil Gas Samples		0		

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WOODARD & CURRAN Engineering • Science • Operations					Geoprobe Log				Boring No.
					Georgia - Pacific				SB-33
41 Hutchins Drive					Old Town, Maine				
Portland, ME 04102									
START DATE: May 25, 2006					BORING DEPTH: 10'				
CONTRACTOR: Environmental Services Network					TEMPORARY WELL DEPTH: 10'				
DRILLER: David Dionne					TEMPORARY RISER LENGTH: N/A				
GEOLOGIST: Danielle Taylor					TEMPORARY SCREEN LENGTH: 4'				
DRILLING METHOD: Direct Push					SLOT SIZE: N/A				
Depth (feet)	**Sample Range**	**Sample Number**	**Rec. Length**	**PID (ppm)**	**Field Description and Remarks**				
0	0-4		2.5'		0 to 0.5	Asphalt			
					0.5 to 1.5	Medium to fine brown SAND, some coarse, black			
						staining, strong odor			
					1.5 to 2.5	Brown coarse SAND			
5						End of soil investigation at 4' bgs			
					* sheen, floating debris and strong odor in water sample (6-10' bgs)				
10									
15									
20									
Groundwater Level Data									
Date	Time	Groundwater Depth (ft)			Summary				
						Soil Samples		0	
						Groundwater Samples		1	
						Soil Gas Samples		0	

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WOODARD & CURRAN Engineering • Science • Operations					Geoprobe Log				Boring No.
					Georgia Pacific				SB-35
41 Hutchins Drive					Old Town, Maine				
Portland, ME 04102									
START DATE: May 25, 2006					BORING DEPTH: 6"'				
CONTRACTOR:					TEMPORARY WELL DEPTH:				N/A
DRILLER:					TEMPORARY RISER LENGTH: N/A				
GEOLOGIST: Josh Bowe					TEMPORARY SCREEN LENGTH: N/A				
DRILLING METHOD: Hand-Dug Soil Sample					SLOT SIZE: N/A				
Depth (feet)	**Sample Range**	**Sample Number**	**Rec. Length**	**PID (ppm)**	**Field Description and Remarks**				
0	0-6"				0-6"	f/m brown sands, some brown silt.			
						Some coarse gravel			
						Hand-dug to ~ 6"			
5									
10									
15									
20									
Groundwater Level Data									
Date	Time	Groundwater Depth (ft)			Summary				
						Soil Samples		1	
						Groundwater Samples		0	
						Soil Gas Samples		0	

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					Geoprobe Log				**Boring No.**
					Georgia - Pacific				**SB-37**
41 Hutchins Drive					**Old Town, Maine**				
Portland, ME 04102									
START DATE: May 25, 2006					BORING DEPTH: 11.7'				
CONTRACTOR: Environmental Services Network					TEMPORARY WELL DEPTH: 11.7'				
DRILLER: David Dionne					TEMPORARY RISER LENGTH: N/A				
GEOLOGIST: Danielle Taylor					TEMPORARY SCREEN LENGTH: 5'				
DRILLING METHOD: Direct Push					SLOT SIZE: N/A				
Depth	**Sample**	**Sample**	**Rec.**	**PID**					
(feet)	**Range**	**Number**	**Length**	**(ppm)**	**Field Description and Remarks**				
0	0-4		3.5		0 to 0.5	Brown coarse to medium SAND, some black staining			
					1 to 3.5	Brown silt and clay			
	4-8		3		4 to 7	Brown silt and clay			
5									
						End of soil investigation at 8' bgs			
10									
15									
20									

Groundwater Level Data									
Date	Time	Groundwater Depth (ft)			Summary				
						Soil Samples		1	
						Groundwater Samples		2	
						Soil Gas Samples		0	

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					Geoprobe Log					**Boring No.**
					Georgia - Pacific					**SB-38**
41 Hutchins Drive					**Old Town, Maine**					
Portland, ME 04102										
START DATE: May 25, 2006						BORING DEPTH: 12'				
CONTRACTOR: Environmental Services Network						TEMPORARY WELL DEPTH: 12'				
DRILLER: David Dionne						TEMPORARY RISER LENGTH: N/A				
GEOLOGIST: Danielle Taylor						TEMPORARY SCREEN LENGTH: 4'				
DRILLING METHOD: Direct Push						SLOT SIZE: N/A				
Depth (feet)	**Sample Range**	**Sample Number**	**Rec. Length**	**PID (ppm)**	**Field Description and Remarks**					
0	0-4		3.5		0 to 1	Brown coarse sand, some black staining				
					1 to 3.5	Brown fine to very fine SAND, some silt and clay				
	4-8		3		4 to 7	Brown silt and clay				
5										
	8-12		3		8 to 11	Brown silt and clay, trace brown coarse SAND				
10										
						End of investigation at 12' bgs				
15										
20										
Groundwater Level Data										
Date	Time	Groundwater Depth (ft)			Summary					
						Soil Samples		1		
						Groundwater Samples		1		
						Soil Gas Samples		0		

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					Geoprobe Log				**Boring No.**
					Georgia - Pacific				**SB-39**
41 Hutchins Drive					**Old Town, Maine**				
Portland, ME 04102									
START DATE: May 25, 2006					BORING DEPTH: 16'				
CONTRACTOR: Environmental Services Network					TEMPORARY WELL DEPTH: 16'				
DRILLER: David Dionne					TEMPORARY RISER LENGTH: N/A				
GEOLOGIST: Danielle Taylor					TEMPORARY SCREEN LENGTH: N/A				
DRILLING METHOD: Direct Push					SLOT SIZE: N/A				
Depth	**Sample**	**Sample**	**Rec.**	**PID**					
(feet)	**Range**	**Number**	**Length**	**(ppm)**	**Field Description and Remarks**				
0	0-4		3.5		0 to 0.5	Asphalt, Brown coarse to medium SAND			
					0.5 to 3	Brown clay and silt, some black staining			
	4-8		3		4 to 7	Brown silt and clay			
5									
	8-12		4		8 to 12	Brown silt and clay			
10									
	12-16		4		12 to 16	Brown silt and clay, dry			
15									
20									
Groundwater Level Data									
Date	Time	Groundwater Depth (ft)			Summary				
						Soil Samples		0	
						Groundwater Samples		0	
						Soil Gas Samples		0	

Appendix B
Geoprobe Logs
Georgia-Pacific Paper Mill
Old Town, Maine

WOODARD & CURRAN
Engineering • Science • Operations

						Geoprobe Log				**Boring No.**
						Georgia - Pacific				**SB-40**
41 Hutchins Drive						**Old Town, Maine**				
Portland, ME 04102										
START DATE: May 25, 2006						BORING DEPTH: 5.5'				
CONTRACTOR: Environmental Services Network						TEMPORARY WELL DEPTH: 5.5'				
DRILLER: David Dionne						TEMPORARY RISER LENGTH: N/A				
GEOLOGIST: Danielle Taylor						TEMPORARY SCREEN LENGTH: N/A				
DRILLING METHOD: Direct Push						SLOT SIZE: N/A				
Depth	**Sample**	**Sample**	**Rec.**	**PID**						
(feet)	**Range**	**Number**	**Length**	**(ppm)**	**Field Description and Remarks**					
0	0-4		3		0 to 0.5	Brown fine to medium uniform SAND, some brown				
						silt				
					0.5 to 1.5	Black wood-like material, organics				
					1.5 to 3	Light brown fine SAND and silt				
	4-5.5		1.5		4 to 5.5	Brown medium to coarse SAND to fine light brown				
5						SAND, some silt				
						weathered bedrock in drive shoe				
						Refusal at 5.5 bgs				
10										
15										
20										
Groundwater Level Data										
Date	Time	Groundwater Depth (ft)				Summary				
						Soil Samples		0		
						Groundwater Samples		1		
						Soil Gas Samples		0		

Appendix B
Geoprobe Logs
Georgia-Pacific Paper Mill
Old Town, Maine

WOODARD & CURRAN
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					Geoprobe Log				**Boring No.**
					Georgia - Pacific				**SB-41**
41 Hutchins Drive					**Old Town, Maine**				
Portland, ME 04102									
START DATE: May 25, 2006					BORING DEPTH: 6'				
CONTRACTOR: Environmental Services Network					TEMPORARY WELL DEPTH: 6'				
DRILLER: David Dionne					TEMPORARY RISER LENGTH: N/A				
GEOLOGIST: Danielle Taylor					TEMPORARY SCREEN LENGTH: 4'				
DRILLING METHOD: Direct Push					SLOT SIZE: N/A				
Depth	**Sample**	**Sample**	**Rec.**	**PID**					
(feet)	**Range**	**Number**	**Length**	**(ppm)**	**Field Description and Remarks**				
0	0-4		3		0 to 1	Organics, brown fine to medium SAND, trace coarse			
						SAND			
					1 to 3	Light brown trace orange silt and fine SAND			
	4-6		2		4 to 5	Brown coarse to fine SAND, some silt			
5					5 to 6	Brown fine SAND and silt, some medium SAND			
						till in drive shoe			
						Refusal at 6' bgs			
10									
15									
20									
Groundwater Level Data									
Date	Time	Groundwater Depth (ft)			Summary				
						Soil Samples		1	
						Groundwater Samples		0	
						Soil Gas Samples		0	

Appendix B
Monitoring Well Boring Log
Georgia-Pacific Paper Mill
Old Town, Maine

WOODARD & CURRAN
Engineering • Science • Operations

BORING LOG

41 Hutchins Drive
Portland, Maine

SITE LOCATION

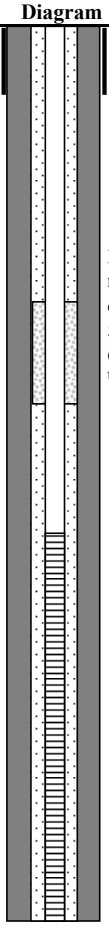
Property: Georgia Pacific Old Town Mill
Address: 24 Portland Street, Old Town, Maine

BORING/ WELL No.
MW-01

PAGE
1 of 1

START DATE	5/26/2006	BORING DEPTH	18	feet bgs	
CONTRACTOR	New Hampshire Borings	WELL DEPTH	18	feet bgs	
DRILLER	Steve Garside	RISER LENGTH	7.5	feet	
GEOLOGIST	Jason House	SCREEN LENGTH	10	feet	
DRILLING METHOD	HSA	SLOT SIZE	0.01	inches	
BORING TYPE	4.25" Radius ID	G.WATER DEPTH	7.52 pvc	feet bgs	

Depth	Sample No.	Blows per 6 inches	Rec. Length	PID (ppm)	Field Description and Remarks	Well Construction Diagram
	SS1		2'		0.0 to 0.7' Topsoil, very fine to fine dark brown SAND, some silt	Road Box
1					0.7 to 1.3' Brown very fine SAND, silt, some coarse SAND	
					1.3 to 2.0' Coarse SAND and GRAVEL	
2						
						Native backfill
3						
4	SS2		3.3'		4 to 4.3' Coarse SAND and GRAVEL, some orange fine material	
					4.3 to 4.5' Orange/brown very fine SAND and silt, trace coarse SAND	Bentonite chips 5.5 to 6.5 ft. bgs
5						
					4.5 to 6.5' Brown very fine to fine SAND, trace coarse SAND, some brown silt	
6						
					6.5 to 7.3' Weathered rock/cobble	Top of sand pack (6.5 ft. bgs)
7						
8	SS3		4.0'		8.0 to 10.0 Brown very fine to fine SAND and silt, trace medium to coarse SAND	
9						
10					10.0 to 12' Brown very fine to fine SAND and silt, trace medium to coarse SAND	Top of Screen (7.77
11						
12	SS4		1.0'		12 to 13.0' Brown very fine to fine SAND, some round-subangular medium to coarse SAND	
13						
					End of soil investigation@ 13' bgs	
14						Bottom



Appendix B
Monitoring Well Boring Log
Georgia-Pacific Paper Mill
Old Town, Maine

WOODARD & CURRAN
Engineering • Science • Operations

BORING LOG

41 Hutchins Drive
Portland, Maine

SITE LOCATION

Property: Georgia Pacific Old Town Mill
Address: 24 Portland Street, Old Town, Maine

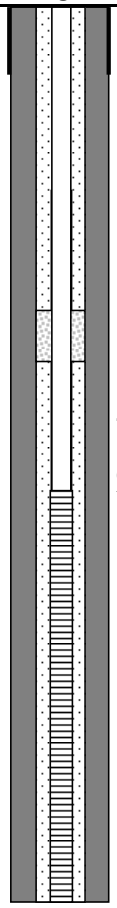
BORING/ WELL No.
MW-02

PAGE
1 of 2

START DATE	5/26/2006	BORING DEPTH	23.5	feet bgs	
CONTRACTOR	New Hampshire Borings	WELL DEPTH	23.5	feet bgs	
DRILLER	Ken	RISER LENGTH	8.2	feet	
GEOLOGIST	Jason House	SCREEN LENGTH	15	feet	
DRILLING METHOD	Drive and Wash	SLOT SIZE	0.01	inches	
BORING TYPE	4" ID Flush Joint Casing	G.WATER DEPTH	~14.5	feet bgs	

Depth	Sample No.	Blows per 6 inches	Rec. Length	PID (ppm)	Field Description and Remarks
	SS-1	8	0.7'		0.0 to 0.2' Fine SANDY LOAM, some pebbles
1		12			0.2 to 0.7' Brown fine to coarse SAND, dry, loose
		15			
2		20			
3					
4					
5	SS-2	15	0.5'		Brown to gray, medium to coarse SAND, some silt, some clay, moist,
		8			loose to medium dense.
6		7			
		12			
7					
8					
9					
10	SS-3	12	0.9'		10 to 10.5' Gray to olive gray, fine to medium SAND, some silt, clay,
		8			and pebbles, medium dense.
11		8			10.5 to 10.9' Black ash material, trace brick fragments, loose. FILL
		21			
12					
13					
14					
15	SS-4	6	0.1		Gray, coarse sand, brick and concrete fragments, loose, wet. FILL

Well Construction Diagram



Appendix B
Monitoring Well Boring Log
Georgia-Pacific Paper Mill
Old Town, Maine

WOODARD & CURRAN
Engineering • Science • Operations

BORING LOG

41 Hutchins Drive
Portland, Maine

SITE LOCATION

Property: Georgia Pacific Old Town Mill
Address: 24 Portland Street, Old Town, Maine

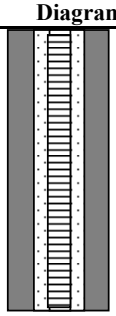
BORING/ WELL No.
MW-02

PAGE
2 of 2

START DATE	5/26/2006	BORING DEPTH	23.5	feet bgs
CONTRACTOR	NH Borings	WELL DEPTH	23.5	feet bgs
DRILLER	Ken	RISER LENGTH	8	feet
GEOLOGIST	Jason House	SCREEN LENGTH	15	feet
DRILLING METHOD	Drive and Wash	SLOT SIZE	0.01	inches
BORING TYPE	4"ID Flush Joint Casing	G.WATER DEPTH	~14	feet bgs

Depth	Sample No.	Blows per 6 inches	Rec. Length	PID (ppm)	Field Description and Remarks
20	SS-5	12	0.9		Black, concrete fragments, red brick fragments, medium sand wet, loose.
		14			
21		14			
		13			
22					
23					
24					
25					
26					
27					
28					
29					
30					
31					

Well Construction Diagram



Bottom of Screen 23.17 ft. bgs

Botto m of MW-02,

Appendix B
Monitoring Well Boring Log
Georgia-Pacific Paper Mill
Old Town, Maine

WOODARD & CURRAN
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BORING LOG

41 Hutchins Drive
Portland, Maine

SITE LOCATION

Property: Georgia Pacific Old Town Mill

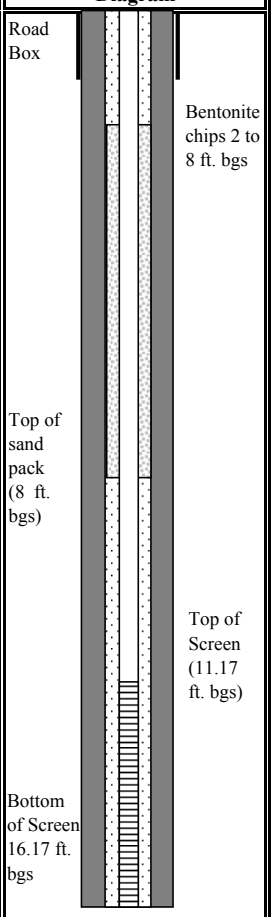
Address: 24 Portland Street, Old Town, Maine

BORING/ WELL No.
MW-03

START DATE	5/25/2006	BORING DEPTH	16.5	feet bgs	
CONTRACTOR	New Hampshire Borings	WELL DEPTH	16.5	feet bgs	**PAGE**
DRILLER	Steve Garside	RISER LENGTH	10	feet	
GEOLOGIST	Jason House	SCREEN LENGTH	5	feet	1 of 1
DRILLING METHOD	HSA	SLOT SIZE	0.01	inches	
BORING TYPE	4.25" Radius ID	G.WATER DEPTH	~11.5	feet bgs	

Depth	Sample No.	Blows per 6 inches	Rec. Length	PID (ppm)	Field Description and Remarks
1					
2					
3					
4	SS-1	4	1.2	0	Fine sand, trace medium sand, some brick fragments, dry, medium dense. FILL
		5			
5		8			
		12			
6					
7					
8					
9	SS-2	12	0.2'	0	Gray, concrete chips. FILL
		9			
10		8			
		5			
11					
12					
13					
14	SS-3	4	1.0'	0	Brown, medium to coarse sand, some brick fragments, black liquid globules, sheen on water from sample, wet, loose.
		3			
15		4			

Well Construction Diagram



Appendix B
Monitoring Well Boring Log
Georgia-Pacific Paper Mill
Old Town, Maine

WOODARD & CURRAN
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BORING LOG

41 Hutchins Drive
Portland, Maine

SITE LOCATION

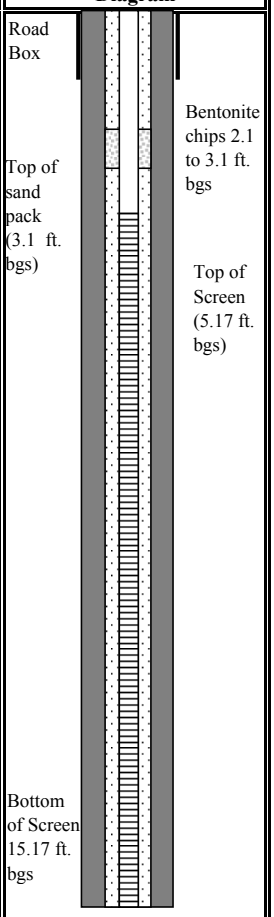
Property: Georgia Pacific Old Town Mill
Address: 24 Portland Street, Old Town, Maine

BORING/ WELL No. MW-04

START DATE	5/25/2006	BORING DEPTH	15.5	feet bgs	
CONTRACTOR	New Hampshire Borings	WELL DEPTH	15.5	feet bgs	**PAGE**
DRILLER	Steve Garside	RISER LENGTH	5	feet	
GEOLOGIST	Jason House	SCREEN LENGTH	10	feet	1 of 1
DRILLING METHOD	HSA	SLOT SIZE	0.01	inches	
BORING TYPE	4.25" Radius ID	G.WATER DEPTH	~6.5	feet bgs	

Depth	Sample No.	Blows per 6 inches	Rec. Length	PID (ppm)	Field Description and Remarks
1					
2					
3					
4	SS-1	13	0.8'	0	Brick fragments with some pebbles and trace medium sand. Loose and dry to moist. FILL
		13			
5		8			
		11			
6					
7					
8					
9	SS-2		1.3'	1.6	9.0 - 9.5' brown to black, medium to fine sand, some silt, wet, loose sulfurous odor.
		1			
10		21			9.5 -10.3' wood.
		41			
11					
12					
13					
14	SS-3	50/0	0.0'		Black phyllite rock fragment logded in spoon tip.
15					

Well Construction Diagram



Appendix B
Monitoring Well Boring Log
Georgia-Pacific Paper Mill
Old Town, Maine

WOODARD & CURRAN
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BORING LOG

41 Hutchins Drive
Portland, Maine

SITE LOCATION

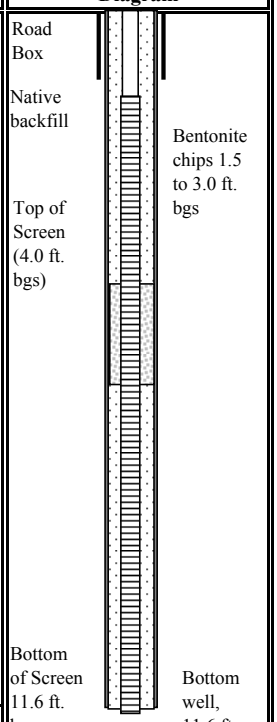
Property: Georgia Pacific Old Town Mill
Address: 24 Portland Street, Old Town, Maine

BORING/ WELL No.
MW-05

START DATE	5/26/2006	BORING DEPTH	11.6	feet bgs	
CONTRACTOR	Environmental Services Network	WELL DEPTH	11.6	feet bgs	**PAGE**
DRILLER	David Dionne	RISER LENGTH	6	feet	
GEOLOGIST	Danielle Taylor	SCREEN LENGTH	7.6	feet	1 of 1
DRILLING METHOD	Direct push	SLOT SIZE	0.01	inches	
BORING TYPE	1.0" PVC	G.WATER DEPTH	~ 4'	feet bgs	

Depth	Sample No.	Blows per 6 inches	Rec. Length	PID (ppm)	Field Description and Remarks
	SS1		2'		0.0 to 0.6' Topsoil, very fine to fine brown SAND
1					0.6 to 1.5' Brown very fine to fine SAND trace coarse SAND
					1.5 to 1.8' Coarse SAND and GRAVEL
2					1.8 to 2.0' Reddish brown fine to medium SAND
3					
4	SS2		3.7'		4 to 4.7' Reddish brown round to subangualar SAND
					4.7 to 5.1' Gray fine to very fine SAND, trace of angular medium SAND, some gray silt
5					
					5.1 to 5.6' Gray/brown very fine to fine SAND, some silt, some coarse SAND
6					
					5.6 to 6.3' Gray/brown very fine to fine SAND, more round/subang coarse SAND
7					
					6.3 to 7.7' Brown medium to fine SAND, some silt
8	SS3		2		8 to 9 Brown very fine sand, brown silt
					9 to 10 Brown medium to coarse SAND, some coarse gravel
9					
10					10.0 to 12' Brown very fine to fine SAND and silt, trace medium to coarse SAND
11					
					Refusal at 11.6 bgs
12	SS4				
13					
14					
15					

Well Construction Diagram



Appendix B
Test Pit Log
Georgia-Pacific Paper Mill
Old Town, Maine

WOODARD & CURRAN Engineering • Science • Operations **TEST PIT LOG** 41 Hutchins Drive Portland, Maine 04096	**SITE LOCATION** GP Old Town, Maine	**TEST PIT No.** TP-101

JOB NUMBER	218736	START DATE	05/30/06	PAGE
CONTRACTOR	Mike Thorton	FINISH DATE	05/30/06	1 of 1
EQUIPMENT	Excavator	TOTAL DEPTH (FT.)	3'	
INSPECTOR	Steve Bartlett	G. WATER DEPTH (FT.)	not measured	
		DIMENSION (L x W)	10' x 2.5'	

Depth (feet)	Field Description and Remarks
0	0-0.5': brown sandy LOAM
1	
2	
	0.5-3': cobbles, brick, logs, cm sand
3	Bottom of test pit at 3 feet; bedrock
4	
5	
6	
7	
8	
9	
10	

NOTES:
Depth to groundwater was not measured because groundwater was not encountered in test pit.
Coordinates: 19 T 0528335, 4971695 NAD 27

Appendix B
Test Pit Log
Georgia-Pacific Paper Mill
Old Town, Maine

WOODARD & CURRAN
Engineering • Science • Operations

TEST PIT LOG

41 Hutchins Drive
Portland, Maine 04096

SITE LOCATION

GP

Old Town, Maine

TEST PIT No.

TP-102

JOB NUMBER	218736	START DATE	05/30/06	PAGE
CONTRACTOR	Mike Thorton	FINISH DATE	05/30/06	1 of 1
EQUIPMENT	Excavator	TOTAL DEPTH (FT.)	3'	
INSPECTOR	Steve Bartlett	G. WATER DEPTH (FT.)	not measured	
		DIMENSION (L x W)	10' x 2.5'	

Depth (feet)	Field Description and Remarks
0	0-0.5': brown sandy LOAM
	0.5-3': fill material, cobbles, asphalt, trace/little sludge
1	
2	
3	Bottom of test pit at 3 feet
4	
5	
6	
7	
8	
9	
10	

NOTES:
Depth to groundwater was not measured because groundwater was not encountered in test pit.
Coordinates: 19 T 0528309, 4971692

Appendix B
Test Pit Log
Georgia-Pacific Paper Mill
Old Town, Maine

WOODARD & CURRAN
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TEST PIT LOG

41 Hutchins Drive
Portland, Maine 04096

SITE LOCATION

GP

Old Town, Maine

TEST PIT No.

TP-103

JOB NUMBER	218736	START DATE	05/30/06	PAGE
CONTRACTOR	Mike Thorton	FINISH DATE	05/30/06	1 of 1
EQUIPMENT	Excavator	TOTAL DEPTH (FT.)	0.5'	
INSPECTOR	Steve Bartlett	G. WATER DEPTH (FT.)	not measured	
		DIMENSION (L x W)	10' x 2.5'	

Depth (feet)	Field Description and Remarks
0	0-0.5': grey sandy fill material, bedrock at 0.5'
	Bottom of test pit at 0.5 feet
1	
2	
3	
4	
5	
6	
7	
8	
9	
10	

NOTES:
Depth to groundwater was not measured because groundwater was not encountered in test pit.
Coordinates: 19 T 0528285, 4971692

Appendix B
Test Pit Log
Georgia-Pacific Paper Mill
Old Town, Maine

WOODARD & CURRAN
Engineering • Science • Operations

TEST PIT LOG

41 Hutchins Drive
Portland, Maine 04096

SITE LOCATION

GP

Old Town, Maine

TEST PIT No.

TP-104

JOB NUMBER	218736	START DATE	05/30/06	PAGE
CONTRACTOR	Mike Thorton	FINISH DATE	05/30/06	1 of 1
EQUIPMENT	Excavator	TOTAL DEPTH (FT.)	3'	
INSPECTOR	Steve Bartlett	G. WATER DEPTH (FT.)	not measured	
		DIMENSION (L x W)	10' x 2.5'	

Depth (feet)	Field Description and Remarks
0	0-3': grey, sandy fill material
1	
2	
3	Bottom of test pit at 3 feet
4	
5	
6	
7	
8	
9	
10	

NOTES:
Depth to groundwater was not measured because groundwater was not encountered in test pit.
Coordinates: 19 T 0528262, 4971662

Appendix B
Test Pit Log
Georgia-Pacific Paper Mill
Old Town, Maine

WOODARD & CURRAN
Engineering • Science • Operations

TEST PIT LOG

41 Hutchins Drive
Portland, Maine 04096

SITE LOCATION

GP

Old Town, Maine

TEST PIT No.

TP-105

JOB NUMBER	218736	START DATE	05/30/06	PAGE
CONTRACTOR	Mike Thorton	FINISH DATE	05/30/06	1 of 1
EQUIPMENT	Excavator	TOTAL DEPTH (FT.)	9'	
INSPECTOR	Steve Bartlett	G. WATER DEPTH (FT.)	not measured	
		DIMENSION (L x W)	10' x 2.5'	

Depth (feet)	Field Description and Remarks
0	0-8': grey, sludge interbedded with wood chips
1	
2	
3	
4	
5	
6	
7	
8	8-9': light grey, fine clay (wet)
9	Bottom of test pit at 9 feet.
10	

NOTES:
Depth to groundwater was not measured because groundwater was not encountered in test pit.
Coordinates: 19T 0528329, 4971695
Composite soil sample (SS01) collected from TP-105; VOCs, SVOCs, RCRA8 metals

Appendix B
Test Pit Log
Georgia-Pacific Paper Mill
Old Town, Maine

WOODARD & CURRAN Engineering • Science • Operations **TEST PIT LOG** 41 Hutchins Drive Portland, Maine 04096	**SITE LOCATION** GP Old Town, Maine	**TEST PIT No.** TP-106

JOB NUMBER	218736	START DATE	05/30/06	PAGE
CONTRACTOR	Mike Thorton	FINISH DATE	05/30/06	1 of 1
EQUIPMENT	Excavator	TOTAL DEPTH (FT.)	4'	
INSPECTOR	Steve Bartlett	G. WATER DEPTH (FT.)	not measured	
		DIMENSION (L x W)	10' x 2.5'	

Depth (feet)	Field Description and Remarks
0	0-0.5': brown sandy loam
1	
2	
3	
4	0.5-4': brown sand, gravel, cobbles, some plastic bags
	Bottom of test pit at 4 feet.
5	
6	
7	
8	
9	
10	

NOTES:
Depth to groundwater was not measured because groundwater was not encountered in test pit.
Coordinates: 19T 0528244, 4971614

Appendix B
Test Pit Log
Georgia-Pacific Paper Mill
Old Town, Maine

WOODARD & CURRAN Engineering • Science • Operations **TEST PIT LOG** 41 Hutchins Drive Portland, Maine 04096	**SITE LOCATION** GP Old Town, Maine	**TEST PIT No.** TP-107

JOB NUMBER	218736	START DATE	05/30/06	PAGE
CONTRACTOR	Mike Thorton	FINISH DATE	05/30/06	1 of 1
EQUIPMENT	Excavator	TOTAL DEPTH (FT.)	5'	
INSPECTOR	Steve Bartlett	G. WATER DEPTH (FT.)	4'	
		DIMENSION (L x W)	10' x 2.5'	

Depth (feet)	Field Description and Remarks
0	0-0.5': brown sandy loam
1	
2	
3	0.5-3': brown sand, gravel, cobbles
4	4': water table
5	3-5': plastic sheets, wood, metal debris, drum tops, fill, and chips
	Bottom of test pit at 5 feet
6	
7	
8	
9	
10	

NOTES:
Composite soil sample (SS02) collected from TP-105 for SVOCs. Weathered fuel oil smell coming from test pit.
Coordinates: 19T 0528218, 4971592

Appendix B
Test Pit Log
Georgia-Pacific Paper Mill
Old Town, Maine

WOODARD & CURRAN Engineering • Science • Operations **TEST PIT LOG** 41 Hutchins Drive Portland, Maine 04096	**SITE LOCATION** GP Old Town, Maine	**TEST PIT No.** TP-108

JOB NUMBER	218736	START DATE	05/30/06	PAGE
CONTRACTOR	Mike Thorton	FINISH DATE	05/30/06	1 of 1
EQUIPMENT	Excavator	TOTAL DEPTH (FT.)	3'	
INSPECTOR	Steve Bartlett	G. WATER DEPTH (FT.)	not measured	
		DIMENSION (L x W)	10' x 2.5'	

Depth (feet)	Field Description and Remarks
0	0-0.5': brown sandy loam
1	
2	
3	0.5-3': garbage, plastic, metal coils, drum top unmarked, painted wood debris, did not dig deeper because of coils
	Bottom of test pit at 3 feet
4	
5	
6	
7	
8	
9	
10	

NOTES:
Depth to groundwater was not measured because groundwater was not encountered in test pit.
Coordinates: 19T 0528217, 4971560

Appendix B
Test Pit Log
Georgia-Pacific Paper Mill
Old Town, Maine

WOODARD & CURRAN Engineering • Science • Operations **TEST PIT LOG** 41 Hutchins Drive Portland, Maine 04096	**SITE LOCATION** GP Old Town, Maine	**TEST PIT No.** TP-109

JOB NUMBER	218736	START DATE	05/30/06	PAGE
CONTRACTOR	Mike Thorton	FINISH DATE	05/30/06	1 of 1
EQUIPMENT	Excavator	TOTAL DEPTH (FT.)	5'	
INSPECTOR	Steve Bartlett	G. WATER DEPTH (FT.)	not measured	
		DIMENSION (L x W)	10' x 2.5'	

Depth (feet)	Field Description and Remarks
0	0-0.5': brown sandy loam
1	
2	
3	
4	
5	0.5-5': anthropogenic fill and soil, conveyor belt piece, metal, drum top
	Bottom of test pit at 5 feet
6	
7	
8	
9	
10	

NOTES:
Depth to groundwater was not measured because groundwater was not encountered in test pit.
Coordinates: 19T 0528232, 4971560

Appendix B
Test Pit Log
Georgia-Pacific Paper Mill
Old Town, Maine

WOODARD & CURRAN Engineering • Science • Operations **TEST PIT LOG** 41 Hutchins Drive Portland, Maine 04096	**SITE LOCATION** GP Old Town, Maine	**TEST PIT No.** TP-110

JOB NUMBER	218736	START DATE	05/30/06	PAGE
CONTRACTOR	Mike Thorton	FINISH DATE	05/30/06	1 of 1
EQUIPMENT	Excavator	TOTAL DEPTH (FT.)	3'	
INSPECTOR	Steve Bartlett	G. WATER DEPTH (FT.)	not measured	
		DIMENSION (L x W)	10' x 2.5'	

Depth (feet)	Field Description and Remarks
0	0-0.5': brown sandy loam
1	
2	
3	0.5-3': anthropogenic fill, cobbles, soil, 1" wire cable
	Bottom of test pit at 3 feet
4	
5	
6	
7	
8	
9	
10	

NOTES:
Depth to groundwater was not measured because groundwater was not encountered in test pit.
Coordinates: 19T 0520255, 4971593

Appendix B
Test Pit Log
Georgia-Pacific Paper Mill
Old Town, Maine

WOODARD & CURRAN Engineering • Science • Operations **TEST PIT LOG** 41 Hutchins Drive Portland, Maine 04096	**SITE LOCATION** GP Old Town, Maine	**TEST PIT No.** TP-111

JOB NUMBER	218736	START DATE	05/30/06	PAGE
CONTRACTOR	Mike Thorton	FINISH DATE	05/30/06	1 of 1
EQUIPMENT	Excavator	TOTAL DEPTH (FT.)	4'	
INSPECTOR	Steve Bartlett	G. WATER DEPTH (FT.)	4'	
		DIMENSION (L x W)	10' x 2.5'	

Depth (feet)	Field Description and Remarks
0	0-0.5': brown sandy loam
1	
2	
3	
4	0.5-4': anthropogenic fill, wod debris, plastic gasket fitting, cardboard boxes, sheet metal
	Bottom of test pit at 4 feet
5	
6	
7	
8	
9	
10	

NOTES:
Coordinates: 19T 0528276, 4971564

Appendix B
Test Pit Log
Georgia-Pacific Paper Mill
Old Town, Maine

WOODARD & CURRAN Engineering • Science • Operations **TEST PIT LOG** 41 Hutchins Drive Portland, Maine 04096	**SITE LOCATION** GP Old Town, Maine	**TEST PIT No.** TP-112

JOB NUMBER	218736	START DATE	05/30/06	PAGE
CONTRACTOR	Mike Thorton	FINISH DATE	05/30/06	1 of 1
EQUIPMENT	Excavator	TOTAL DEPTH (FT.)	3'	
INSPECTOR	Steve Bartlett	G. WATER DEPTH (FT.)	not measured	
		DIMENSION (L x W)	10' x 2.5'	

Depth (feet)	Field Description and Remarks
0	
1	0-1': grey sandy fill
2	
3	1-3': sand/cobbles; bedrock at 3 feet
	Bottom of test pit at 3 feet
4	
5	
6	
7	
8	
9	
10	

NOTES:
Coordinates: 19T 0523290, 4971604
Depth to groundwater was not measured because groundwater was not encountered in test pit.

Appendix B
Test Pit Log
Georgia-Pacific Paper Mill
Old Town, Maine

WOODARD & CURRAN Engineering • Science • Operations **TEST PIT LOG** 41 Hutchins Drive Portland, Maine 04096	**SITE LOCATION** GP Old Town, Maine	**TEST PIT No.** TP-113

JOB NUMBER	218736	START DATE	05/30/06	PAGE
CONTRACTOR	Mike Thorton	FINISH DATE	05/30/06	1 of 1
EQUIPMENT	Excavator	TOTAL DEPTH (FT.)	3'	
INSPECTOR	Steve Bartlett	G. WATER DEPTH (FT.)	not measured	
		DIMENSION (L x W)	10' x 2.5'	

Depth (feet)	Field Description and Remarks
0	
1	0-1': wood chips
2	
3	
4	1-4': fine grey silt with cobbles and boulders
	Bottom of test pit at 3 feet
5	
6	
7	
8	
9	
10	

NOTES:
Coordinates: 19T 0528297, 4971716
Depth to groundwater was not measured because groundwater was not encountered in test pit.

Appendix B
Test Pit Log
Georgia-Pacific Paper Mill
Old Town, Maine

WOODARD & CURRAN Engineering • Science • Operations **TEST PIT LOG** 41 Hutchins Drive Portland, Maine 04096	**SITE LOCATION** GP Old Town, Maine	**TEST PIT No.** TP-114

JOB NUMBER	218736	START DATE	05/30/06	PAGE
CONTRACTOR	Mike Thorton	FINISH DATE	05/30/06	1 of 1
EQUIPMENT	Excavator	TOTAL DEPTH (FT.)	0.5'	
INSPECTOR	Steve Bartlett	G. WATER DEPTH (FT.)	not measured	
		DIMENSION (L x W)	10' x 2.5'	

Depth (feet)	Field Description and Remarks
0	
	0-0.5': wood chips, bedrock at 0.5'
1	Bottom of test pit at 0.5 feet
2	
3	
4	
5	
6	
7	
8	
9	
10	

NOTES:
Coordinates: 19T 0528222, 4971731
Depth to groundwater was not measured because groundwater was not encountered in test pit.

Appendix B
Test Pit Log
Georgia-Pacific Paper Mill
Old Town, Maine

WOODARD & CURRAN Engineering • Science • Operations **TEST PIT LOG** 41 Hutchins Drive Portland, Maine 04096	**SITE LOCATION** GP Old Town, Maine	**TEST PIT No.** TP-115

JOB NUMBER	218736	START DATE	05/30/06	PAGE
CONTRACTOR	Mike Thorton	FINISH DATE	05/30/06	1 of 1
EQUIPMENT	Excavator	TOTAL DEPTH (FT.)	4'	
INSPECTOR	Steve Bartlett	G. WATER DEPTH (FT.)	not measured	
		DIMENSION (L x W)	10' x 2.5'	

Depth (feet)	Field Description and Remarks
0	
1	
2	
3	
4	0-4': fill soil, cobbles, boulders
	Bottom of test pit at 4 feet
5	
6	
7	
8	
9	
10	

NOTES:
Coordinates: 19T 0528149, 4971739
Depth to groundwater was not measured because groundwater was not encountered in test pit.

Appendix B
Test Pit Log
Georgia-Pacific Paper Mill
Old Town, Maine

WOODARD & CURRAN Engineering • Science • Operations **TEST PIT LOG** 41 Hutchins Drive Portland, Maine 04096	**SITE LOCATION** GP Old Town, Maine	**TEST PIT No.** TP-116

JOB NUMBER	218736	START DATE	05/30/06	PAGE
CONTRACTOR	Mike Thorton	FINISH DATE	05/30/06	1 of 1
EQUIPMENT	Excavator	TOTAL DEPTH (FT.)	4'	
INSPECTOR	Steve Bartlett	G. WATER DEPTH (FT.)	not measured	
		DIMENSION (L x W)	10' x 2.5'	

Depth (feet)	Field Description and Remarks
0	
1	
2	0-2': sludge
3	
4	2-4': brown sand, cobbles Bottom of test pit at 4 feet
5	
6	
7	
8	
9	
10	

NOTES:
Coordinates: 19T 0528182, 4971729
Depth to groundwater was not measured because groundwater was not encountered in test pit.
Test pit in edge of sludge landfill ramp.
SS03 collected in 0-2' interval

Appendix B
Test Pit Log
Georgia-Pacific Paper Mill
Old Town, Maine

WOODARD & CURRAN Engineering • Science • Operations **TEST PIT LOG** 41 Hutchins Drive Portland, Maine 04096	**SITE LOCATION** GP Old Town, Maine	**TEST PIT No.** TP-117

JOB NUMBER	218736	START DATE	05/30/06	PAGE
CONTRACTOR	Mike Thorton	FINISH DATE	05/30/06	1 of 1
EQUIPMENT	Excavator	TOTAL DEPTH (FT.)	4'	
INSPECTOR	Steve Bartlett	G. WATER DEPTH (FT.)	not measured	
		DIMENSION (L x W)	10' x 2.5'	

Depth (feet)	Field Description and Remarks
0	
1	
2	0-2': sludge
3	
4	2-4': brown sand, cobbles
	Bottom of test pit at 4 feet
5	
6	
7	
8	
9	
10	

NOTES:
Coordinates: 19T 0528185, 4971820
Depth to groundwater was not measured because groundwater was not encountered in test pit.
Test pit in edge of sludge landfill pile.

Appendix B
Test Pit Log
Georgia-Pacific Paper Mill
Old Town, Maine

WOODARD & CURRAN Engineering • Science • Operations **TEST PIT LOG** 41 Hutchins Drive Portland, Maine 04096	**SITE LOCATION** GP Old Town, Maine	**TEST PIT No.** TP-118

JOB NUMBER	218736	START DATE	05/30/06	PAGE
CONTRACTOR	Mike Thorton	FINISH DATE	05/30/06	1 of 1
EQUIPMENT	Excavator	TOTAL DEPTH (FT.)	2'	
INSPECTOR	Steve Bartlett	G. WATER DEPTH (FT.)	not measured	
		DIMENSION (L x W)	10' x 2.5'	

Depth (feet)	Field Description and Remarks
0	
1	
2	0-2': wood chips, bedrock at 2'
	Bottom of test pit at 2 feet
3	
4	
5	
6	
7	
8	
9	
10	

NOTES:
Coordinates: 19T 0528290, 4971875
Depth to groundwater was not measured because groundwater was not encountered in test pit.

Appendix B
Test Pit Log
Georgia-Pacific Paper Mill
Old Town, Maine

WOODARD & CURRAN Engineering • Science • Operations **TEST PIT LOG** 41 Hutchins Drive Portland, Maine 04096	**SITE LOCATION** GP Old Town, Maine	**TEST PIT No.** TP-119

JOB NUMBER	218736	START DATE	05/30/06	PAGE
CONTRACTOR	Mike Thorton	FINISH DATE	05/30/06	1 of 1
EQUIPMENT	Excavator	TOTAL DEPTH (FT.)	2'	
INSPECTOR	Steve Bartlett	G. WATER DEPTH (FT.)	not measured	
		DIMENSION (L x W)	10' x 2.5'	

Depth (feet)	Field Description and Remarks
0	
1	
2	
3	
4	0-4': grey ash-like material interbedded with wood chips
	4-4.5': brown mc sand
5	Bottom of test pit at 4.5 feet
6	
7	
8	
9	
10	

NOTES:
Coordinates: 19T 0528290, 4971875
Depth to groundwater was not measured because groundwater was not encountered in test pit.
Composite soil sample (SS04) collected from TP-119 for SVOCs, VOCs, and RCRA8 metals.

Appendix B
Test Pit Log
Georgia-Pacific Paper Mill
Old Town, Maine

WOODARD & CURRAN Engineering • Science • Operations **TEST PIT LOG** 41 Hutchins Drive Portland, Maine 04096	**SITE LOCATION** GP Old Town, Maine	**TEST PIT No.** TP-120

JOB NUMBER	218736	START DATE	05/30/06	PAGE
CONTRACTOR	Mike Thorton	FINISH DATE	05/30/06	1 of 1
EQUIPMENT	Excavator	TOTAL DEPTH (FT.)	4'	
INSPECTOR	Steve Bartlett	G. WATER DEPTH (FT.)	not measured	
		DIMENSION (L x W)	10' x 2.5'	

Depth (feet)	Field Description and Remarks
0	0-0.5': wood chips
1	
2	
3	
4	0.5-4': brown sandy fill, cobbles, gravel
	Bottom of test pit at 4 feet
5	
6	
7	
8	
9	
10	

NOTES:
Coordinates: 19T 0528346, 4971787
Depth to groundwater was not measured because groundwater was not encountered in test pit.

Appendix B
Test Pit Log
Georgia-Pacific Paper Mill
Old Town, Maine

WOODARD & CURRAN Engineering • Science • Operations **TEST PIT LOG** 41 Hutchins Drive Portland, Maine 04096	**SITE LOCATION** GP Old Town, Maine	**TEST PIT No.** TP-121

JOB NUMBER	218736	START DATE	05/30/06	PAGE
CONTRACTOR	Mike Thorton	FINISH DATE	05/30/06	1 of 1
EQUIPMENT	Excavator	TOTAL DEPTH (FT.)	4'	
INSPECTOR	Steve Bartlett	G. WATER DEPTH (FT.)	not measured	
		DIMENSION (L x W)	10' x 2.5'	

Depth (feet)	Field Description and Remarks
0	
1	
2	
3	
4	0-4': grey ash-like material
5	4-5': brown cmf sand and gravel
	Bottom of test pit at 5 feet
6	
7	
8	
9	
10	

NOTES:
Coordinates: 19T 0528345, 4971758
Depth to groundwater was not measured because groundwater was not encountered in test pit.
Test pit in the side of pile.

Appendix B
Test Pit Log
Georgia-Pacific Paper Mill
Old Town, Maine

WOODARD & CURRAN Engineering • Science • Operations **TEST PIT LOG** 41 Hutchins Drive Portland, Maine 04096	**SITE LOCATION** GP Old Town, Maine	**TEST PIT No.** TP-122

JOB NUMBER	218736	START DATE	05/30/06	PAGE
CONTRACTOR	Mike Thorton	FINISH DATE	05/30/06	1 of 1
EQUIPMENT	Excavator	TOTAL DEPTH (FT.)	4'	
INSPECTOR	Steve Bartlett	G. WATER DEPTH (FT.)	not measured	
		DIMENSION (L x W)	10' x 2.5'	

Depth (feet)	Field Description and Remarks
0	
1	
2	
3	
4	0-4': grey sludge, interbedded with wood chips
	Bottom of test pit at 5 feet
5	
6	
7	
8	
9	
10	

NOTES:
Coordinates: 19T 0528345, 4971758
Depth to groundwater was not measured because groundwater was not encountered in test pit.
Test pit on top of sludge landfill heap.
Composite soil sample (SS05) collected from TP-122 for SVOCs, VOCs, and RCRA8 metals.

APPENDIX C: MOBILE LABORATORY RESULTS

ESN
Environmental Services Network
North Atlantic
Direct Push Sampling
In-Situ Remediation
Mobile & Fixed Base Chemical Analysis
Integrated Site Characterization

Certificate of Analysis

Sample Identification	Method Blank
Depth (ft)	NA
Sample Matrix	soil
Units:	ug/kg
% Solid	100.0%
Dilution	1
Sample Preservation	4 deg C.
Sample Volume	2 oz.
Date Sampled	5/25/2006
Date Analyzed	5/25/2006

Client:	Woodard & Curran, Inc.
Site:	GP Mill
Loc.:	Old Town, Maine
Project No.:	
Chemist:	D. M. Twomey

Analytical Method: USEPA SW-846 8260B
Instrument: Agilent 5890 GCD MS
Detectors: Mass spectrometer
Column: DB-5, 30 meter, .25 um
Conc. System: Tekmar 3000 Purge & Trap System
Trap: Vocarb 3000

Method:
USEPA SW846 METHOD 8260B

Compound	RL ug/kg	Results	Compound	RL ug/kg	Results	Compound	RL ug/kg	Results
Chloromethane	50	<	1,2-Dichloroethane	50	<	Styrene	50	<
Dichlorodifluoromethane	50	<	1,1-Dichloropropene	50	<	trans-1,4-Dichloro-2-butene	50	<
Chloroethane	50	<	Carbon Tetrachloride	50	<	cis-1,4,-Dichloro-2-butene	50	<
Vinyl chloride	50	<	Benzene	50	<	1,1,2,2-Tetrachloroethane	50	<
Bromomethane	100	<	2-Chloroethanol	250	<	1,2,3-Trichloropropane	50	<
Trichlorofluoromethane	50	<	1,2-Dichloropropane	50	<	Isopropylbenzene	50	<
Diethyl ether	250	<	Trichloroethene	50	<	Bromobenzene	50	<
Acetonitrile	250	<	Bromodichloromethane	50	<	2-Chlorotoluene	100	<
Idomethane	250	<	Dibromomethane	50	<	4-Chlorotoluene	100	<
1,1-Dichloroethene	50	<	Methyl methacrylate	250	<	n-propylbenzene	50	<
Carbon Disulfide	100	<	1,4-Dioxane	25	<	1,3,5-Trimethylbenzene	100	<
Allyl chloride	250	<	cis-1,3-Dichloropropene	50	<	Pentachloroethane	50	<
Methylene chloride	500	<	trans-1,3-Dichloropropene	50	<	1,2,4-Trimethylbenzene	100	<
Acrylonitrile	250	<	2-Nitropropane	250	<	tert-butylbenzene	50	<
1,1,2 - Trichlorotrifluoroethane	50	<	Toluene	50	<	sec-butylbenzene	50	<
trans-1,2-Dichloroethene	50	<	1,1,2-Trichloroethane	50	<	1,3-Dichlorobenzene	100	<
MTBE	50	<	Ethylmethacrylate	250	<	1,4-Dichlorobenzene	100	<
Chloroform	50	<	1,3-Dichloropropane	50	<	p-iso-Propyltoluene	50	<
1,1-Dichloroethane	50	<	Dibromochloromethane	50	<	1,2-Dichlorobenzene	100	<
Chloroprene	100	<	1,2-Dibromomethane	50	<	n-butylbenzene	50	<
2,2-Dichloropropane	50	<	Tetrachloroethene	50	<	1,2-Bromo-3-chloropropane	50	<
Methyl acrylate	250	<	Chlorobenzene	50	<	Nitrobenzene	250	<
cis-1,2,-Dichloroethene	50	<	1,1,1,2-Tetrachloroethane	50	<	1,2,4-Trichlorobenzene	50	<
Tetrahydrofuran	250	<	Ethylbenzene	50	<	Naphthalene	25	<
Methyacrylonitrile	250	<	m/p-xylene	50	<	1,2,3-Trichlorobenzene	50	<
Isobutyl alcohol	250	<	Bromoform	100	<	Hexachlorobutadiene	50	<
1,1,1-Trichloroethane	50	<	o-xylene	50	<			

Legend:
< = less than reporting limits
J = less than reporting limit, above detection levels
E = estimated, above calibration range

Surrogate Recovery	%
Dibromofluoromethane	98
Toluene-d8	90
4-Bromofluorobenzene	90

Notes:

Chemist/Reviewer: RMB

Date: 5/31/06

PO Box 6752
Fax: 207-883-1912
15 Holly St., Ste 102
Scarborough, ME 04070
800-880-0533
info@esnnorth.com

ESN North Atlantic
Environmental Services Network
Direct Push Sampling In-Situ Remediation
Mobile & Fixed Base Chemical Analysis Integrated Site Characterization

Certificate of Analysis

Sample Identification	Method Blank
Depth (ft)	NA
Sample Matrix	water
Units:	ug/L
% Solid	NA
Dilution	1
Sample Preservation	4 deg C.
Sample Volume	40 mL.
Date Sampled	5/24/2006
Date Analyzed	5/24/2006

Client:	Woodard & Curran, Inc.
Site:	GP Mill
Loc.:	Old Town, Maine
Project No.:	
Chemist:	D. M. Twomey

Analytical Method: USEPA SW-846 8260B
Instrument: Agilent 5890 GCD MS
Detectors: Mass spectrometer
Column: DB-5, 30 meter, .25 um
Conc. System: Tekmar 3000 Purge & Trap System
Trap: Vocarb 3000

Method:
USEPA SW846 METHOD 8260B

Compound	RL ug/L	Results	Compound	RL ug/L	Results	Compound	RL ug/L	Results
Chloromethane	1	<	1,2-Dichloroethane	1	<	Styrene	1	<
Dichlorodifluoromethane	1	<	1,1-Dichloropropene	1	<	trans-1,4-Dichloro-2-butene	1	<
Chloroethane	1	<	Carbon Tetrachloride	1	<	cis-1,4,-Dichloro-2-butene	1	<
Vinyl chloride	1	<	Benzene	1	<	1,1,2,2-Tetrachloroethane	1	<
Bromomethane	2	<	2-Chloroethanol	5	<	1,2,3-Trichloropropane	1	<
Trichlorofluoromethane	1	<	1,2-Dichloropropane	1	<	Isopropylbenzene	1	<
Diethyl ether	5	<	Trichloroethene	1	<	Bromobenzene	1	<
Acetonitrile	5	<	Bromodichloromethane	1	<	2-Chlorotoluene	2	<
Idomethane	5	<	Dibromomethane	1	<	4-Chlorotoluene	2	<
1,1-Dichloroethene	1	<	Methyl methacrylate	5	<	n-propylbenzene	1	<
Carbon Disulfide	2	<	1,4-Dioxane	25	<	1,3,5-Trimethylbenzene	2	<
Allyl chloride	5	<	cis-1,3-Dichloropropene	1	<	Pentachloroethane	1	<
Methylene chloride	10	<	trans-1,3-Dichloropropene	1	<	1,2,4-Trimethylbenzene	2	<
Acrylonitrile	5	<	2-Nitropropane	5	<	tert-butylbenzene	1	<
1,1,2 - Trichlorotrifluoroethane	1	<	Toluene	1	<	sec-butylbenzene	1	<
trans-1,2-Dichloroethene	1	<	1,1,2-Trichloroethane	1	<	1,3-Dichlorobenzene	2	<
MTBE	1	<	Ethylmethacrylate	5	<	1,4-Dichlorobenzene	2	<
Chloroform	1	<	1,3-Dichloropropane	1	<	p-iso-Propyltoluene	1	<
1,1-Dichloroethane	1	<	Dibromochloromethane	1	<	1,2-Dichlorobenzene	2	<
Chloroprene	2	<	1,2-Dibromomethane	1	<	n-butylbenzene	1	<
2,2-Dichloropropane	1	<	Tetrachloroethene	1	<	1,2-Bromo-3-chloropropane	1	<
Methyl acrylate	5	<	Chlorobenzene	1	<	Nitrobenzene	5	<
cis-1,2,-Dichloroethene	1	<	1,1,1,2-Tetrachloroethane	1	<	1,2,4-Trichlorobenzene	1	<
Tetrahydrofuran	5	<	Ethylbenzene	1	<	Naphthalene	25	<
Methyacrylonitrile	5	<	m/p-xylene	1	<	1,2,3-Trichlorobenzene	1	<
Isobutyl alcohol	5	<	Bromoform	2	<	Hexachlorobutadiene	1	<
1,1,1-Trichloroethane	1	<	o-xylene	1	<			

Legend:
< = less than reporting limits
J = less than reporting limit, above detection levels
E = estimated, above calibration range

Surrogate Recovery	%
Dibromofluoromethane	101
Toluene-d8	96
4-Bromofluorobenzene	95

Notes:

Chemist/Reviewer: RMB

Date: 5/31/06

PO Box 6752
Fax: 207-883-1912
15 Holly St., Ste 102
Scarborough, ME 04070
800-880-0533
info@esnnorth.com

ESN North Atlantic
Environmental Services Network
Direct Push Sampling | In-Situ Remediation | Mobile & Fixed Base Chemical Analysis | Integrated Site Characterization

Certificate of Analysis

Sample Identification	SB-1
Depth (ft)	6-10'
Sample Matrix	water
Units:	ug/L
% Solid	NA
Dilution	1
Sample Preservation	4 deg C.
Sample Volume	40 mL.
Date Sampled	5/25/2006
Date Analyzed	5/25/2006

Client:	**Woodard & Curran, Inc.**
Site:	**GP Mill**
Loc.:	**Old Town, Maine**
Project No.:	
Chemist:	D. M. Twomey

Analytical Method: USEPA SW-846 8260B
Instrument: Agilent 5890 GCD MS
Detectors: Mass spectrometer
Column: DB-5, 30 meter, .25 um
Conc. System: Tekmar 3000 Purge & Trap System
Trap: Vocarb 3000

Method:
USEPA SW846 METHOD 8260B

Compound	RL ug/L	Results	Compound	RL ug/L	Results	Compound	RL ug/L	Results
Chloromethane	1	<	1,2-Dichloroethane	1	<	Styrene	1	<
Dichlorodifluoromethane	1	<	1,1-Dichloropropene	1	<	trans-1,4-Dichloro-2-butene	1	<
Chloroethane	1	<	Carbon Tetrachloride	1	<	cis-1,4,-Dichloro-2-butene	1	<
Vinyl chloride	1	<	Benzene	1	<	1,1,2,2-Tetrachloroethane	1	<
Bromomethane	2	<	2-Chloroethanol	5	<	1,2,3-Trichloropropane	1	<
Trichlorofluoromethane	1	<	1,2-Dichloropropane	1	<	Isopropylbenzene	1	<
Diethyl ether	5	<	Trichloroethene	1	<	Bromobenzene	1	<
Acetonitrile	5	<	Bromodichloromethane	1	<	2-Chlorotoluene	2	<
Idomethane	5	<	Dibromomethane	1	<	4-Chlorotoluene	2	<
1,1-Dichloroethene	1	<	Methyl methacrylate	5	<	n-propylbenzene	1	<
Carbon Disulfide	2	<	1,4-Dioxane	25	<	1,3,5-Trimethylbenzene	2	<
Allyl chloride	5	<	cis-1,3-Dichloropropene	1	<	Pentachloroethane	1	<
Methylene chloride	10	<	trans-1,3-Dichloropropene	1	<	1,2,4-Trimethylbenzene	2	<
Acrylonitrile	5	<	2-Nitropropane	5	<	tert-butylbenzene	1	<
1,1,2 - Trichlorotrifluoroethane	1	<	Toluene	1	<	sec-butylbenzene	1	<
trans-1,2-Dichloroethene	1	<	1,1,2-Trichloroethane	1	<	1,3-Dichlorobenzene	2	<
MTBE	1	<	Ethylmethacrylate	5	<	1,4-Dichlorobenzene	2	<
Chloroform	1	<	1,3-Dichloropropane	1	<	p-iso-Propyltoluene	1	<
1,1-Dichloroethane	1	<	Dibromochloromethane	1	<	1,2-Dichlorobenzene	2	<
Chloroprene	2	<	1,2-Dibromomethane	1	<	n-butylbenzene	1	<
2,2-Dichloropropane	1	<	Tetrachloroethene	1	<	1,2-Bromo-3-chloropropane	1	<
Methyl acrylate	5	<	Chlorobenzene	1	<	Nitrobenzene	5	<
cis-1,2,-Dichloroethene	1	<	1,1,1,2-Tetrachloroethane	1	<	1,2,4-Trichlorobenzene	1	<
Tetrahydrofuran	5	<	Ethylbenzene	1	<	Naphthalene	25	<
Methyacrylonitrile	5	<	m/p-xylene	1	<	1,2,3-Trichlorobenzene	1	<
Isobutyl alcohol	5	<	Bromoform	2	<	Hexachlorobutadiene	1	<
1,1,1-Trichloroethane	1	<	o-xylene	1	<			

Legend:
< = less than reporting limits
J = less than reporting limit, above detection levels
E = estimated, above calibration range

Surrogate Recovery	%
Dibromofluoromethane	97
Toluene-d8	109
4-Bromofluorobenzene	92

Notes:

Chemist/Reviewer: RMB

Date: 5/31/06

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Certificate of Analysis

Sample Identification	SB-1
Depth (ft)	0-4'
Sample Matrix	soil
Units:	ug/kg
% Solid	88.9%
Dilution	1
Sample Preservation	4 deg C.
Sample Volume	2 oz.
Date Sampled	5/25/2006
Date Analyzed	5/25/2006

Client:	**Woodard & Curran, Inc.**
Site:	**GP Mill**
Loc.:	**Old Town, Maine**
Project No.:	
Chemist:	D. M. Twomey

Analytical Method: USEPA SW-846 8260B
Instrument: Agilent 5890 GCD MS
Detectors: Mass spectrometer
Column: DB-5, 30 meter, .25 um
Conc. System: Tekmar 3000 Purge & Trap System
Trap: Vocarb 3000

Method:
USEPA SW846 METHOD 8260B

Compound	RL ug/kg	Results	Compound	RL ug/kg	Results	Compound	RL ug/kg	Results
Chloromethane	50	<	1,2-Dichloroethane	50	<	Styrene	50	<
Dichlorodifluoromethane	50	<	1,1-Dichloropropene	50	<	trans-1,4-Dichloro-2-butene	50	<
Chloroethane	50	<	Carbon Tetrachloride	50	<	cis-1,4,-Dichloro-2-butene	50	<
Vinyl chloride	50	<	Benzene	50	<	1,1,2,2-Tetrachloroethane	50	<
Bromomethane	100	<	2-Chloroethanol	250	<	1,2,3-Trichloropropane	50	<
Trichlorofluoromethane	50	<	1,2-Dichloropropane	50	<	Isopropylbenzene	50	<
Diethyl ether	250	<	Trichloroethene	50	<	Bromobenzene	50	<
Acetonitrile	250	<	Bromodichloromethane	50	<	2-Chlorotoluene	100	<
Idomethane	250	<	Dibromomethane	50	<	4-Chlorotoluene	100	<
1,1-Dichloroethene	50	<	Methyl methacrylate	250	<	n-propylbenzene	50	<
Carbon Disulfide	100	<	1,4-Dioxane	25	<	1,3,5-Trimethylbenzene	100	<
Allyl chloride	250	<	cis-1,3-Dichloropropene	50	<	Pentachloroethane	50	<
Methylene chloride	500	<	trans-1,3-Dichloropropene	50	<	1,2,4-Trimethylbenzene	100	<
Acrylonitrile	250	<	2-Nitropropane	250	<	tert-butylbenzene	50	<
1,1,2 - Trichlorotrifluoroethane	50	<	Toluene	50	<	sec-butylbenzene	50	<
trans-1,2-Dichloroethene	50	<	1,1,2-Trichloroethane	50	<	1,3-Dichlorobenzene	100	<
MTBE	50	<	Ethylmethacrylate	250	<	1,4-Dichlorobenzene	100	<
Chloroform	50	<	1,3-Dichloropropane	50	<	p-iso-Propyltoluene	50	<
1,1-Dichloroethane	50	<	Dibromochloromethane	50	<	1,2-Dichlorobenzene	100	<
Chloroprene	100	<	1,2-Dibromomethane	50	<	n-butylbenzene	50	<
2,2-Dichloropropane	50	<	Tetrachloroethene	50	<	1,2-Bromo-3-chloropropane	50	<
Methyl acrylate	250	<	Chlorobenzene	50	<	Nitrobenzene	250	<
cis-1,2,-Dichloroethene	50	<	1,1,1,2-Tetrachloroethane	50	<	1,2,4-Trichlorobenzene	50	<
Tetrahydrofuran	250	<	Ethylbenzene	50	<	Naphthalene	25	<
Methyacrylonitrile	250	<	m/p-xylene	50	<	1,2,3-Trichlorobenzene	50	<
Isobutyl alcohol	250	<	Bromoform	100	<	Hexachlorobutadiene	50	<
1,1,1-Trichloroethane	50	<	o-xylene	50	<			

Legend:
< = less than reporting limits
J = less than reporting limit, above detection levels
E = estimated, above calibration range

Surrogate Recovery	%
Dibromofluoromethane	79
Toluene-d8	97
4-Bromofluorobenzene	85

Notes:

Chemist/Reviewer: RMB

Date: 5/31/06

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Certificate of Analysis

Sample Identification	SB-18
Depth (ft)	10-14'
Sample Matrix	water
Units:	ug/L
% Solid	NA
Dilution	1
Sample Preservation	4 deg C.
Sample Volume	40 mL.
Date Sampled	5/25/2006
Date Analyzed	5/25/2006

Client:	Woodard & Curran, Inc.
Site:	GP Mill
Loc.:	Old Town, Maine
Project No.:	
Chemist:	D. M. Twomey

Analytical Method: USEPA SW-846 8260B
Instrument: Agilent 5890 GCD MS
Detectors: Mass spectrometer
Column: DB-5, 30 meter, .25 um
Conc. System: Tekmar 3000 Purge & Trap System
Trap: Vocarb 3000

Method:
USEPA SW846 METHOD 8260B

Compound	RL ug/L	Results	Compound	RL ug/L	Results	Compound	RL ug/L	Results
Chloromethane	1	<	1,2-Dichloroethane	1	<	Styrene	1	<
Dichlorodifluoromethane	1	<	1,1-Dichloropropene	1	<	trans-1,4-Dichloro-2-butene	1	<
Chloroethane	1	<	Carbon Tetrachloride	1	<	cis-1,4,-Dichloro-2-butene	1	<
Vinyl chloride	1	<	Benzene	1	<	1,1,2,2-Tetrachloroethane	1	<
Bromomethane	2	<	2-Chloroethanol	5	<	1,2,3-Trichloropropane	1	<
Trichlorofluoromethane	1	<	1,2-Dichloropropane	1	<	Isopropylbenzene	1	<
Diethyl ether	5	<	Trichloroethene	1	<	Bromobenzene	1	<
Acetonitrile	5	<	Bromodichloromethane	1	<	2-Chlorotoluene	2	<
Idomethane	5	<	Dibromomethane	1	<	4-Chlorotoluene	2	<
1,1-Dichloroethene	1	<	Methyl methacrylate	5	<	n-propylbenzene	1	<
Carbon Disulfide	2	<	1,4-Dioxane	25	<	1,3,5-Trimethylbenzene	2	<
Allyl chloride	5	<	cis-1,3-Dichloropropene	1	<	Pentachloroethane	1	<
Methylene chloride	10	<	trans-1,3-Dichloropropene	1	<	1,2,4-Trimethylbenzene	2	<
Acrylonitrile	5	<	2-Nitropropane	5	<	tert-butylbenzene	1	<
1,1,2 - Trichlorotrifluoroethane	1	<	Toluene	1	<	sec-butylbenzene	1	<
trans-1,2-Dichloroethene	1	<	1,1,2-Trichloroethane	1	<	1,3-Dichlorobenzene	2	<
MTBE	1	<	Ethylmethacrylate	5	<	1,4-Dichlorobenzene	2	<
Chloroform	1	<	1,3-Dichloropropane	1	<	p-iso-Propyltoluene	1	<
1,1-Dichloroethane	1	<	Dibromochloromethane	1	<	1,2-Dichlorobenzene	2	<
Chloroprene	2	<	1,2-Dibromomethane	1	<	n-butylbenzene	1	<
2,2-Dichloropropane	1	<	Tetrachloroethene	1	<	1,2-Bromo-3-chloropropane	1	<
Methyl acrylate	5	<	Chlorobenzene	1	<	Nitrobenzene	5	<
cis-1,2,-Dichloroethene	1	<	1,1,1,2-Tetrachloroethane	1	<	1,2,4-Trichlorobenzene	1	<
Tetrahydrofuran	5	<	Ethylbenzene	1	<	Naphthalene	25	<
Methyacrylonitrile	5	<	m/p-xylene	1	<	1,2,3-Trichlorobenzene	1	<
Isobutyl alcohol	5	<	Bromoform	2	<	Hexachlorobutadiene	1	<
1,1,1-Trichloroethane	1	<	o-xylene	1	<			

Legend:
< = less than reporting limits
J = less than reporting limit, above detection levels
E = estimated, above calibration range

Surrogate Recovery	%
Dibromofluoromethane	87
Toluene-d8	100
4-Bromofluorobenzene	85

Notes:

Chemist/Reviewer: RMB

Date: 5/31/06

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Certificate of Analysis

Sample Identification	SB-11
Depth (ft)	6-9'
Sample Matrix	water
Units:	ug/L
% Solid	NA
Dilution	1
Sample Preservation	4 deg C.
Sample Volume	40 mL.
Date Sampled	5/24/2006
Date Analyzed	5/24/2006

Client:	**Woodard & Curran, Inc.**
Site:	**GP Mill**
Loc.:	**Old Town, Maine**
Project No.:	
Chemist:	D. M. Twomey

Analytical Method: USEPA SW-846 8260B
Instrument: Agilent 5890 GCD MS
Detectors: Mass spectrometer
Column: DB-5, 30 meter, .25 um
Conc. System: Tekmar 3000 Purge & Trap System
Trap: Vocarb 3000

Method:
USEPA SW846 METHOD 8260B

Compound	RL ug/L	Results	Compound	RL ug/L	Results	Compound	RL ug/L	Results
Chloromethane	1	<	1,2-Dichloroethane	1	<	Styrene	1	<
Dichlorodifluoromethane	1	<	1,1-Dichloropropene	1	<	trans-1,4-Dichloro-2-butene	1	<
Chloroethane	1	<	Carbon Tetrachloride	1	<	cis-1,4,-Dichloro-2-butene	1	<
Vinyl chloride	1	<	Benzene	1	<	1,1,2,2-Tetrachloroethane	1	<
Bromomethane	2	<	2-Chloroethanol	5	<	1,2,3-Trichloropropane	1	<
Trichlorofluoromethane	1	<	1,2-Dichloropropane	1	<	Isopropylbenzene	1	<
Diethyl ether	5	<	Trichloroethene	1	<	Bromobenzene	1	<
Acetonitrile	5	<	Bromodichloromethane	1	<	2-Chlorotoluene	2	<
Idomethane	5	<	Dibromomethane	1	<	4-Chlorotoluene	2	<
1,1-Dichloroethene	1	<	Methyl methacrylate	5	<	n-propylbenzene	1	<
Carbon Disulfide	2	<	1,4-Dioxane	25	<	1,3,5-Trimethylbenzene	2	<
Allyl chloride	5	<	cis-1,3-Dichloropropene	1	<	Pentachloroethane	1	<
Methylene chloride	10	<	trans-1,3-Dichloropropene	1	<	1,2,4-Trimethylbenzene	2	<
Acrylonitrile	5	<	2-Nitropropane	5	<	tert-butylbenzene	1	<
1,1,2 - Trichlorotrifluoroethane	1	<	Toluene	1	<	sec-butylbenzene	1	<
trans-1,2-Dichloroethene	1	<	1,1,2-Trichloroethane	1	<	1,3-Dichlorobenzene	2	<
MTBE	1	<	Ethylmethacrylate	5	<	1,4-Dichlorobenzene	2	<
Chloroform	1	<	1,3-Dichloropropane	1	<	p-iso-Propyltoluene	1	<
1,1-Dichloroethane	1	<	Dibromochloromethane	1	<	1,2-Dichlorobenzene	2	<
Chloroprene	2	<	1,2-Dibromomethane	1	<	n-butylbenzene	1	<
2,2-Dichloropropane	1	<	Tetrachloroethene	1	<	1,2-Bromo-3-chloropropane	1	<
Methyl acrylate	5	<	Chlorobenzene	1	<	Nitrobenzene	5	<
cis-1,2,-Dichloroethene	1	<	1,1,1,2-Tetrachloroethane	1	<	1,2,4-Trichlorobenzene	1	<
Tetrahydrofuran	5	<	Ethylbenzene	1	<	Naphthalene	25	<
Methyacrylonitrile	5	<	m/p-xylene	1	<	1,2,3-Trichlorobenzene	1	<
Isobutyl alcohol	5	<	Bromoform	2	<	Hexachlorobutadiene	1	<
1,1,1-Trichloroethane	1	<	o-xylene	1	<			

Legend:
< = less than reporting limits
J = less than reporting limit, above detection levels
E = estimated, above calibration range

Surrogate Recovery	%
Dibromofluoromethane	92
Toluene-d8	96
4-Bromofluorobenzene	97

Notes:

Chemist/Reviewer: RMB

Date: 5/31/06

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Certificate of Analysis

Sample Identification	SB-12
Depth (ft)	6-9'
Sample Matrix	water
Units:	ug/L
% Solid	NA
Dilution	1
Sample Preservation	4 deg C.
Sample Volume	40 mL.
Date Sampled	5/24/2006
Date Analyzed	5/24/2006

Client:	**Woodard & Curran, Inc.**
Site:	**GP Mill**
Loc.:	**Old Town, Maine**
Project No.:	
Chemist:	D. M. Twomey

Analytical Method: USEPA SW-846 8260B
Instrument: Agilent 5890 GCD MS
Detectors: Mass spectrometer
Column: DB-5, 30 meter, .25 um
Conc. System: Tekmar 3000 Purge & Trap System
Trap: Vocarb 3000

Method:
USEPA SW846 METHOD 8260B

Compound	RL ug/L	Results	Compound	RL ug/L	Results	Compound	RL ug/L	Results
Chloromethane	1	<	1,2-Dichloroethane	1	<	Styrene	1	<
Dichlorodifluoromethane	1	<	1,1-Dichloropropene	1	<	trans-1,4-Dichloro-2-butene	1	<
Chloroethane	1	<	Carbon Tetrachloride	1	<	cis-1,4,-Dichloro-2-butene	1	<
Vinyl chloride	1	<	Benzene	1	<	1,1,2,2-Tetrachloroethane	1	<
Bromomethane	2	<	2-Chloroethanol	5	<	1,2,3-Trichloropropane	1	<
Trichlorofluoromethane	1	<	1,2-Dichloropropane	1	<	Isopropylbenzene	1	<
Diethyl ether	5	<	Trichloroethene	1	<	Bromobenzene	1	<
Acetonitrile	5	<	Bromodichloromethane	1	<	2-Chlorotoluene	2	<
Idomethane	5	<	Dibromomethane	1	<	4-Chlorotoluene	2	<
1,1-Dichloroethene	1	<	Methyl methacrylate	5	<	n-propylbenzene	1	<
Carbon Disulfide	2	<	1,4-Dioxane	25	<	1,3,5-Trimethylbenzene	2	<
Allyl chloride	5	<	cis-1,3-Dichloropropene	1	<	Pentachloroethane	1	<
Methylene chloride	10	<	trans-1,3-Dichloropropene	1	<	1,2,4-Trimethylbenzene	2	<
Acrylonitrile	5	<	2-Nitropropane	5	<	tert-butylbenzene	1	<
1,1,2 - Trichlorotrifluoroethane	1	<	Toluene	1	<	sec-butylbenzene	1	<
trans-1,2-Dichloroethene	1	<	1,1,2-Trichloroethane	1	<	1,3-Dichlorobenzene	2	<
MTBE	1	<	Ethylmethacrylate	5	<	1,4-Dichlorobenzene	2	<
Chloroform	1	<	1,3-Dichloropropane	1	<	p-iso-Propyltoluene	1	<
1,1-Dichloroethane	1	<	Dibromochloromethane	1	<	1,2-Dichlorobenzene	2	<
Chloroprene	2	<	1,2-Dibromomethane	1	<	n-butylbenzene	1	<
2,2-Dichloropropane	1	<	Tetrachloroethene	1	<	1,2-Bromo-3-chloropropane	1	<
Methyl acrylate	5	<	Chlorobenzene	1	<	Nitrobenzene	5	<
cis-1,2,-Dichloroethene	1	<	1,1,1,2-Tetrachloroethane	1	<	1,2,4-Trichlorobenzene	1	<
Tetrahydrofuran	5	<	Ethylbenzene	1	<	Naphthalene	25	<
Methyacrylonitrile	5	<	m/p-xylene	1	<	1,2,3-Trichlorobenzene	1	<
Isobutyl alcohol	5	<	Bromoform	2	<	Hexachlorobutadiene	1	<
1,1,1-Trichloroethane	1	<	o-xylene	1	<			

Legend:
< = less than reporting limits
J = less than reporting limit, above detection levels
E = estimated, above calibration range

Surrogate Recovery	%
Dibromofluoromethane	92
Toluene-d8	95
4-Bromofluorobenzene	90

Notes:

Chemist/Reviewer: RMB

Date: 5/31/06

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Certificate of Analysis

Sample Identification	SB-13
Depth (ft)	6-10'
Sample Matrix	water
Units:	ug/L
% Solid	NA
Dilution	1
Sample Preservation	4 deg C.
Sample Volume	40 mL.
Date Sampled	5/24/2006
Date Analyzed	5/24/2006

Client:	**Woodard & Curran, Inc.**
Site:	**GP Mill**
Loc.:	**Old Town, Maine**
Project No.:	
Chemist:	D. M. Twomey

Analytical Method: USEPA SW-846 8260B
Instrument: Agilent 5890 GCD MS
Detectors: Mass spectrometer
Column: DB-5, 30 meter, .25 um
Conc. System: Tekmar 3000 Purge & Trap System
Trap: Vocarb 3000

Method:
USEPA SW846 METHOD 8260B

Compound	RL ug/L	Results	Compound	RL ug/L	Results	Compound	RL ug/L	Results
Chloromethane	1	<	1,2-Dichloroethane	1	<	Styrene	1	<
Dichlorodifluoromethane	1	<	1,1-Dichloropropene	1	<	trans-1,4-Dichloro-2-butene	1	<
Chloroethane	1	<	Carbon Tetrachloride	1	<	cis-1,4,-Dichloro-2-butene	1	<
Vinyl chloride	1	<	Benzene	1	<	1,1,2,2-Tetrachloroethane	1	<
Bromomethane	2	<	2-Chloroethanol	5	<	1,2,3-Trichloropropane	1	<
Trichlorofluoromethane	1	<	1,2-Dichloropropane	1	<	Isopropylbenzene	1	<
Diethyl ether	5	<	Trichloroethene	1	<	Bromobenzene	1	<
Acetonitrile	5	<	Bromodichloromethane	1	<	2-Chlorotoluene	2	<
Idomethane	5	<	Dibromomethane	1	<	4-Chlorotoluene	2	<
1,1-Dichloroethene	1	2.2	Methyl methacrylate	5	<	n-propylbenzene	1	<
Carbon Disulfide	2	<	1,4-Dioxane	25	<	1,3,5-Trimethylbenzene	2	<
Allyl chloride	5	<	cis-1,3-Dichloropropene	1	<	Pentachloroethane	1	<
Methylene chloride	10	<	trans-1,3-Dichloropropene	1	<	1,2,4-Trimethylbenzene	2	<
Acrylonitrile	5	<	2-Nitropropane	5	<	tert-butylbenzene	1	<
1,1,2 - Trichlorotrifluoroethane	1	<	Toluene	1	<	sec-butylbenzene	1	<
trans-1,2-Dichloroethene	1	<	1,1,2-Trichloroethane	1	<	1,3-Dichlorobenzene	2	<
MTBE	1	<	Ethylmethacrylate	5	<	1,4-Dichlorobenzene	2	<
Chloroform	1	<	1,3-Dichloropropane	1	<	p-iso-Propyltoluene	1	<
1,1-Dichloroethane	1	40	Dibromochloromethane	1	<	1,2-Dichlorobenzene	2	<
Chloroprene	2	<	1,2-Dibromomethane	1	<	n-butylbenzene	1	<
2,2-Dichloropropane	1	<	Tetrachloroethene	1	<	1,2-Bromo-3-chloropropane	1	<
Methyl acrylate	5	<	Chlorobenzene	1	<	Nitrobenzene	5	<
cis-1,2,-Dichloroethene	1	<	1,1,1,2-Tetrachloroethane	1	<	1,2,4-Trichlorobenzene	1	<
Tetrahydrofuran	5	<	Ethylbenzene	1	<	Naphthalene	25	<
Methyacrylonitrile	5	<	m/p-xylene	1	<	1,2,3-Trichlorobenzene	1	<
Isobutyl alcohol	5	<	Bromoform	2	<	Hexachlorobutadiene	1	<
1,1,1-Trichloroethane	1	20	o-xylene	1	<			

Legend:
< = less than reporting limits
J = less than reporting limit, above detection levels
E = estimated, above calibration range

Surrogate Recovery	%
Dibromofluoromethane	89
Toluene-d8	91
4-Bromofluorobenzene	81

Notes:

Chemist/Reviewer: RMB

Date: 5/31/06

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Sample Identification	SB-14
Depth (ft)	6-10'
Sample Matrix	water
Units:	ug/L
% Solid	NA
Dilution	1
Sample Preservation	4 deg C.
Sample Volume	40 mL.
Date Sampled	5/24/2006
Date Analyzed	5/24/2006

Client:	**Woodard & Curran, Inc.**
Site:	**GP Mill**
Loc.:	**Old Town, Maine**
Project No.:	
Chemist:	D. M. Twomey

Analytical Method: USEPA SW-846 8260B
Instrument: Agilent 5890 GCD MS
Detectors: Mass spectrometer
Column: DB-5, 30 meter, .25 um
Conc. System: Tekmar 3000 Purge & Trap System
Trap: Vocarb 3000

Method:
USEPA SW846 METHOD 8260B

Compound	RL ug/L	Results	Compound	RL ug/L	Results	Compound	RL ug/L	Results
Chloromethane	1	<	1,2-Dichloroethane	1	<	Styrene	1	<
Dichlorodifluoromethane	1	<	1,1-Dichloropropene	1	<	trans-1,4-Dichloro-2-butene	1	<
Chloroethane	1	<	Carbon Tetrachloride	1	<	cis-1,4,-Dichloro-2-butene	1	<
Vinyl chloride	1	<	Benzene	1	<	1,1,2,2-Tetrachloroethane	1	<
Bromomethane	2	<	2-Chloroethanol	5	<	1,2,3-Trichloropropane	1	<
Trichlorofluoromethane	1	<	1,2-Dichloropropane	1	<	Isopropylbenzene	1	<
Diethyl ether	5	<	Trichloroethene	1	<	Bromobenzene	1	<
Acetonitrile	5	<	Bromodichloromethane	1	<	2-Chlorotoluene	2	<
Idomethane	5	<	Dibromomethane	1	<	4-Chlorotoluene	2	<
1,1-Dichloroethene	1	<	Methyl methacrylate	5	<	n-propylbenzene	1	<
Carbon Disulfide	2	<	1,4-Dioxane	25	<	1,3,5-Trimethylbenzene	2	<
Allyl chloride	5	<	cis-1,3-Dichloropropene	1	<	Pentachloroethane	1	<
Methylene chloride	10	<	trans-1,3-Dichloropropene	1	<	1,2,4-Trimethylbenzene	2	<
Acrylonitrile	5	<	2-Nitropropane	5	<	tert-butylbenzene	1	<
1,1,2 - Trichlorotrifluoroethane	1	<	Toluene	1	<	sec-butylbenzene	1	<
trans-1,2-Dichloroethene	1	<	1,1,2-Trichloroethane	1	<	1,3-Dichlorobenzene	2	<
MTBE	1	<	Ethylmethacrylate	5	<	1,4-Dichlorobenzene	2	<
Chloroform	1	<	1,3-Dichloropropane	1	<	p-iso-Propyltoluene	1	<
1,1-Dichloroethane	1	<	Dibromochloromethane	1	<	1,2-Dichlorobenzene	2	<
Chloroprene	2	<	1,2-Dibromomethane	1	<	n-butylbenzene	1	<
2,2-Dichloropropane	1	<	Tetrachloroethene	1	<	1,2-Bromo-3-chloropropane	1	<
Methyl acrylate	5	<	Chlorobenzene	1	<	Nitrobenzene	5	<
cis-1,2,-Dichloroethene	1	<	1,1,1,2-Tetrachloroethane	1	<	1,2,4-Trichlorobenzene	1	<
Tetrahydrofuran	5	<	Ethylbenzene	1	<	Naphthalene	25	<
Methyacrylonitrile	5	<	m/p-xylene	1	<	1,2,3-Trichlorobenzene	1	<
Isobutyl alcohol	5	<	Bromoform	2	<	Hexachlorobutadiene	1	<
1,1,1-Trichloroethane	1	<	o-xylene	1	<			

Legend:
< = less than reporting limits
J = less than reporting limit, above detection levels
E = estimated, above calibration range

Surrogate Recovery	%
Dibromofluoromethane	92
Toluene-d8	95
4-Bromofluorobenzene	94

Notes:

Chemist/Reviewer: RMB

Date: 5/31/06

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Certificate of Analysis

Sample Identification	SB-17
Depth (ft)	3.5 - 7.5'
Sample Matrix	water
Units:	ug/L
% Solid	NA
Dilution	1
Sample Preservation	4 deg C.
Sample Volume	40 mL.
Date Sampled	5/24/2006
Date Analyzed	5/24/2006

Client:	**Woodard & Curran, Inc.**
Site:	**GP Mill**
Loc.:	**Old Town, Maine**
Project No.:	
Chemist:	D. M. Twomey

Analytical Method: USEPA SW-846 8260B
Instrument: Agilent 5890 GCD MS
Detectors: Mass spectrometer
Column: DB-5, 30 meter, .25 um
Conc. System: Tekmar 3000 Purge & Trap System
Trap: Vocarb 3000

Method:
USEPA SW846 METHOD 8260B

Compound	RL ug/L	Results	Compound	RL ug/L	Results	Compound	RL ug/L	Results
Chloromethane	1	<	1,2-Dichloroethane	1	<	Styrene	1	<
Dichlorodifluoromethane	1	<	1,1-Dichloropropene	1	<	trans-1,4-Dichloro-2-butene	1	<
Chloroethane	1	<	Carbon Tetrachloride	1	<	cis-1,4,-Dichloro-2-butene	1	<
Vinyl chloride	1	<	Benzene	1	<	1,1,2,2-Tetrachloroethane	1	<
Bromomethane	2	<	2-Chloroethanol	5	<	1,2,3-Trichloropropane	1	<
Trichlorofluoromethane	1	<	1,2-Dichloropropane	1	<	Isopropylbenzene	1	<
Diethyl ether	5	<	Trichloroethene	1	<	Bromobenzene	1	<
Acetonitrile	5	<	Bromodichloromethane	1	<	2-Chlorotoluene	2	<
Idomethane	5	<	Dibromomethane	1	<	4-Chlorotoluene	2	<
1,1-Dichloroethene	1	<	Methyl methacrylate	5	<	n-propylbenzene	1	<
Carbon Disulfide	2	<	1,4-Dioxane	25	<	1,3,5-Trimethylbenzene	2	<
Allyl chloride	5	<	cis-1,3-Dichloropropene	1	<	Pentachloroethane	1	<
Methylene chloride	10	<	trans-1,3-Dichloropropene	1	<	1,2,4-Trimethylbenzene	2	<
Acrylonitrile	5	<	2-Nitropropane	5	<	tert-butylbenzene	1	<
1,1,2 - Trichlorotrifluoroethane	1	<	Toluene	1	<	sec-butylbenzene	1	<
trans-1,2-Dichloroethene	1	<	1,1,2-Trichloroethane	1	<	1,3-Dichlorobenzene	2	<
MTBE	1	<	Ethylmethacrylate	5	<	1,4-Dichlorobenzene	2	<
Chloroform	1	<	1,3-Dichloropropane	1	<	p-iso-Propyltoluene	1	<
1,1-Dichloroethane	1	<	Dibromochloromethane	1	<	1,2-Dichlorobenzene	2	<
Chloroprene	2	<	1,2-Dibromomethane	1	<	n-butylbenzene	1	<
2,2-Dichloropropane	1	<	Tetrachloroethene	1	<	1,2-Bromo-3-chloropropane	1	<
Methyl acrylate	5	<	Chlorobenzene	1	<	Nitrobenzene	5	<
cis-1,2,-Dichloroethene	1	<	1,1,1,2-Tetrachloroethane	1	<	1,2,4-Trichlorobenzene	1	<
Tetrahydrofuran	5	<	Ethylbenzene	1	<	Naphthalene	25	<
Methyacrylonitrile	5	<	m/p-xylene	1	<	1,2,3-Trichlorobenzene	1	<
Isobutyl alcohol	5	<	Bromoform	2	<	Hexachlorobutadiene	1	<
1,1,1-Trichloroethane	1	<	o-xylene	1	<			

Legend:
< = less than reporting limits
J = less than reporting limit, above detection levels
E = estimated, above calibration range

Surrogate Recovery	%
Dibromofluoromethane	86
Toluene-d8	96
4-Bromofluorobenzene	92

Notes:

Chemist/Reviewer: RMB

Date: 5/31/06

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Certificate of Analysis

Sample Identification	SB-15
Depth (ft)	6-10'
Sample Matrix	water
Units:	ug/L
% Solid	NA
Dilution	1
Sample Preservation	4 deg C.
Sample Volume	40 mL.
Date Sampled	5/24/2006
Date Analyzed	5/24/2006

Client:	**Woodard & Curran, Inc.**
Site:	**GP Mill**
Loc.:	**Old Town, Maine**
Project No.:	
Chemist:	D. M. Twomey

Analytical Method: USEPA SW-846 8260B
Instrument: Agilent 5890 GCD MS
Detectors: Mass spectrometer
Column: DB-5, 30 meter, .25 um
Conc. System: Tekmar 3000 Purge & Trap System
Trap: Vocarb 3000

Method:
USEPA SW846 METHOD 8260B

Compound	RL ug/L	Results	Compound	RL ug/L	Results	Compound	RL ug/L	Results
Chloromethane	1	<	1,2-Dichloroethane	1	<	Styrene	1	<
Dichlorodifluoromethane	1	<	1,1-Dichloropropene	1	<	trans-1,4-Dichloro-2-butene	1	<
Chloroethane	1	<	Carbon Tetrachloride	1	<	cis-1,4,-Dichloro-2-butene	1	<
Vinyl chloride	1	<	Benzene	1	<	1,1,2,2-Tetrachloroethane	1	<
Bromomethane	2	<	2-Chloroethanol	5	<	1,2,3-Trichloropropane	1	<
Trichlorofluoromethane	1	<	1,2-Dichloropropane	1	<	Isopropylbenzene	1	<
Diethyl ether	5	<	Trichloroethene	1	<	Bromobenzene	1	<
Acetonitrile	5	<	Bromodichloromethane	1	<	2-Chlorotoluene	2	<
Idomethane	5	<	Dibromomethane	1	<	4-Chlorotoluene	2	<
1,1-Dichloroethene	1	<	Methyl methacrylate	5	<	n-propylbenzene	1	<
Carbon Disulfide	2	<	1,4-Dioxane	25	<	1,3,5-Trimethylbenzene	2	<
Allyl chloride	5	<	cis-1,3-Dichloropropene	1	<	Pentachloroethane	1	<
Methylene chloride	10	<	trans-1,3-Dichloropropene	1	<	1,2,4-Trimethylbenzene	2	<
Acrylonitrile	5	<	2-Nitropropane	5	<	tert-butylbenzene	1	<
1,1,2 - Trichlorotrifluoroethane	1	<	Toluene	1	<	sec-butylbenzene	1	<
trans-1,2-Dichloroethene	1	<	1,1,2-Trichloroethane	1	<	1,3-Dichlorobenzene	2	<
MTBE	1	<	Ethylmethacrylate	5	<	1,4-Dichlorobenzene	2	<
Chloroform	1	<	1,3-Dichloropropane	1	<	p-iso-Propyltoluene	1	<
1,1-Dichloroethane	1	<	Dibromochloromethane	1	<	1,2-Dichlorobenzene	2	<
Chloroprene	2	<	1,2-Dibromomethane	1	<	n-butylbenzene	1	<
2,2-Dichloropropane	1	<	Tetrachloroethene	1	<	1,2-Bromo-3-chloropropane	1	<
Methyl acrylate	5	<	Chlorobenzene	1	<	Nitrobenzene	5	<
cis-1,2,-Dichloroethene	1	<	1,1,1,2-Tetrachloroethane	1	<	1,2,4-Trichlorobenzene	1	<
Tetrahydrofuran	5	<	Ethylbenzene	1	<	Naphthalene	25	<
Methyacrylonitrile	5	<	m/p-xylene	1	<	1,2,3-Trichlorobenzene	1	<
Isobutyl alcohol	5	<	Bromoform	2	<	Hexachlorobutadiene	1	<
1,1,1-Trichloroethane	1	<	o-xylene	1	<			

Legend:
< = less than reporting limits
J = less than reporting limit, above detection levels
E = estimated, above calibration range

Surrogate Recovery	%
Dibromofluoromethane	93
Toluene-d8	92
4-Bromofluorobenzene	91

Notes:

Chemist/Reviewer: RMB

Date: 5/31/06

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Certificate of Analysis

Sample Identification	SB-19
Depth (ft)	2-6'
Sample Matrix	water
Units:	ug/L
% Solid	NA
Dilution	1
Sample Preservation	4 deg C.
Sample Volume	40 mL.
Date Sampled	5/24/2006
Date Analyzed	5/24/2006

Client:	**Woodard & Curran, Inc.**
Site:	**GP Mill**
Loc.:	**Old Town, Maine**
Project No.:	
Chemist:	D. M. Twomey

Analytical Method: USEPA SW-846 8260B
Instrument: Agilent 5890 GCD MS
Detectors: Mass spectrometer
Column: DB-5, 30 meter, .25 um
Conc. System: Tekmar 3000 Purge & Trap System
Trap: Vocarb 3000

Method:
USEPA SW846 METHOD 8260B

Compound	RL ug/L	Results	Compound	RL ug/L	Results	Compound	RL ug/L	Results
Chloromethane	1	<	1,2-Dichloroethane	1	<	Styrene	1	<
Dichlorodifluoromethane	1	<	1,1-Dichloropropene	1	<	trans-1,4-Dichloro-2-butene	1	<
Chloroethane	1	<	Carbon Tetrachloride	1	<	cis-1,4,-Dichloro-2-butene	1	<
Vinyl chloride	1	<	Benzene	1	<	1,1,2,2-Tetrachloroethane	1	<
Bromomethane	2	<	2-Chloroethanol	5	<	1,2,3-Trichloropropane	1	<
Trichlorofluoromethane	1	<	1,2-Dichloropropane	1	<	Isopropylbenzene	1	<
Diethyl ether	5	<	Trichloroethene	1	<	Bromobenzene	1	<
Acetonitrile	5	<	Bromodichloromethane	1	<	2-Chlorotoluene	2	<
Idomethane	5	<	Dibromomethane	1	<	4-Chlorotoluene	2	<
1,1-Dichloroethene	1	<	Methyl methacrylate	5	<	n-propylbenzene	1	<
Carbon Disulfide	2	<	1,4-Dioxane	25	<	1,3,5-Trimethylbenzene	2	<
Allyl chloride	5	<	cis-1,3-Dichloropropene	1	<	Pentachloroethane	1	<
Methylene chloride	10	<	trans-1,3-Dichloropropene	1	<	1,2,4-Trimethylbenzene	2	<
Acrylonitrile	5	<	2-Nitropropane	5	<	tert-butylbenzene	1	<
1,1,2 - Trichlorotrifluoroethane	1	<	Toluene	1	<	sec-butylbenzene	1	<
trans-1,2-Dichloroethene	1	<	1,1,2-Trichloroethane	1	<	1,3-Dichlorobenzene	2	<
MTBE	1	<	Ethylmethacrylate	5	<	1,4-Dichlorobenzene	2	<
Chloroform	1	<	1,3-Dichloropropane	1	<	p-iso-Propyltoluene	1	<
1,1-Dichloroethane	1	<	Dibromochloromethane	1	<	1,2-Dichlorobenzene	2	<
Chloroprene	2	<	1,2-Dibromomethane	1	<	n-butylbenzene	1	<
2,2-Dichloropropane	1	<	Tetrachloroethene	1	<	1,2-Bromo-3-chloropropane	1	<
Methyl acrylate	5	<	Chlorobenzene	1	<	Nitrobenzene	5	<
cis-1,2,-Dichloroethene	1	<	1,1,1,2-Tetrachloroethane	1	<	1,2,4-Trichlorobenzene	1	<
Tetrahydrofuran	5	<	Ethylbenzene	1	<	Naphthalene	25	<
Methyacrylonitrile	5	<	m/p-xylene	1	<	1,2,3-Trichlorobenzene	1	<
Isobutyl alcohol	5	<	Bromoform	2	<	Hexachlorobutadiene	1	<
1,1,1-Trichloroethane	1	<	o-xylene	1	<			

Legend:
< = less than reporting limits
J = less than reporting limit, above detection levels
E = estimated, above calibration range

Surrogate Recovery	%
Dibromofluoromethane	91
Toluene-d8	95
4-Bromofluorobenzene	91

Notes:

Chemist/Reviewer: RMB

Date: 5/31/06

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Certificate of Analysis

Sample Identification	SB-10
Depth (ft)	6-10'
Sample Matrix	water
Units:	ug/L
% Solid	NA
Dilution	1
Sample Preservation	4 deg C.
Sample Volume	40 mL.
Date Sampled	5/24/2006
Date Analyzed	5/24/2006

Client:	**Woodard & Curran, Inc.**
Site:	**GP Mill**
Loc.:	**Old Town, Maine**
Project No.:	
Chemist:	D. M. Twomey

Analytical Method: USEPA SW-846 8260B
Instrument: Agilent 5890 GCD MS
Detectors: Mass spectrometer
Column: DB-5, 30 meter, .25 um
Conc. System: Tekmar 3000 Purge & Trap System
Trap: Vocarb 3000

Method:
USEPA SW846 METHOD 8260B

Compound	RL ug/L	Results	Compound	RL ug/L	Results	Compound	RL ug/L	Results
Chloromethane	1	<	1,2-Dichloroethane	1	<	Styrene	1	<
Dichlorodifluoromethane	1	<	1,1-Dichloropropene	1	<	trans-1,4-Dichloro-2-butene	1	<
Chloroethane	1	<	Carbon Tetrachloride	1	<	cis-1,4,-Dichloro-2-butene	1	<
Vinyl chloride	1	<	Benzene	1	<	1,1,2,2-Tetrachloroethane	1	<
Bromomethane	2	<	2-Chloroethanol	5	<	1,2,3-Trichloropropane	1	<
Trichlorofluoromethane	1	<	1,2-Dichloropropane	1	<	Isopropylbenzene	1	<
Diethyl ether	5	<	Trichloroethene	1	<	Bromobenzene	1	<
Acetonitrile	5	<	Bromodichloromethane	1	<	2-Chlorotoluene	2	<
Idomethane	5	<	Dibromomethane	1	<	4-Chlorotoluene	2	<
1,1-Dichloroethene	1	<	Methyl methacrylate	5	<	n-propylbenzene	1	<
Carbon Disulfide	2	<	1,4-Dioxane	25	<	1,3,5-Trimethylbenzene	2	<
Allyl chloride	5	<	cis-1,3-Dichloropropene	1	<	Pentachloroethane	1	<
Methylene chloride	10	<	trans-1,3-Dichloropropene	1	<	1,2,4-Trimethylbenzene	2	<
Acrylonitrile	5	<	2-Nitropropane	5	<	tert-butylbenzene	1	<
1,1,2 - Trichlorotrifluoroethane	1	<	Toluene	1	<	sec-butylbenzene	1	<
trans-1,2-Dichloroethene	1	<	1,1,2-Trichloroethane	1	<	1,3-Dichlorobenzene	2	<
MTBE	1	<	Ethylmethacrylate	5	<	1,4-Dichlorobenzene	2	<
Chloroform	1	<	1,3-Dichloropropane	1	<	p-iso-Propyltoluene	1	<
1,1-Dichloroethane	1	<	Dibromochloromethane	1	<	1,2-Dichlorobenzene	2	<
Chloroprene	2	<	1,2-Dibromomethane	1	<	n-butylbenzene	1	<
2,2-Dichloropropane	1	<	Tetrachloroethene	1	<	1,2-Bromo-3-chloropropane	1	<
Methyl acrylate	5	<	Chlorobenzene	1	<	Nitrobenzene	5	<
cis-1,2,-Dichloroethene	1	<	1,1,1,2-Tetrachloroethane	1	<	1,2,4-Trichlorobenzene	1	<
Tetrahydrofuran	5	<	Ethylbenzene	1	<	Naphthalene	25	<
Methyacrylonitrile	5	<	m/p-xylene	1	<	1,2,3-Trichlorobenzene	1	<
Isobutyl alcohol	5	<	Bromoform	2	<	Hexachlorobutadiene	1	<
1,1,1-Trichloroethane	1	<	o-xylene	1	<			

Legend:
< = less than reporting limits
J = less than reporting limit, above detection levels
E = estimated, above calibration range

Surrogate Recovery	%
Dibromofluoromethane	87
Toluene-d8	101
4-Bromofluorobenzene	91

Notes:

Chemist/Reviewer: RMB

Date: 5/31/06

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Certificate of Analysis

Sample Identification	SB-20
Depth (ft)	10-14'
Sample Matrix	water
Units:	ug/L
% Solid	NA
Dilution	1
Sample Preservation	4 deg C.
Sample Volume	40 mL.
Date Sampled	5/24/2006
Date Analyzed	5/24/2006

Client:	**Woodard & Curran, Inc.**
Site:	**GP Mill**
Loc.:	**Old Town, Maine**
Project No.:	
Chemist:	D. M. Twomey

Analytical Method: USEPA SW-846 8260B
Instrument: Agilent 5890 GCD MS
Detectors: Mass spectrometer
Column: DB-5, 30 meter, .25 um
Conc. System: Tekmar 3000 Purge & Trap System
Trap: Vocarb 3000

Method:
USEPA SW846 METHOD 8260B

Compound	RL ug/L	Results	Compound	RL ug/L	Results	Compound	RL ug/L	Results
Chloromethane	1	<	1,2-Dichloroethane	1	<	Styrene	1	<
Dichlorodifluoromethane	1	<	1,1-Dichloropropene	1	<	trans-1,4-Dichloro-2-butene	1	<
Chloroethane	1	<	Carbon Tetrachloride	1	<	cis-1,4,-Dichloro-2-butene	1	<
Vinyl chloride	1	<	Benzene	1	<	1,1,2,2-Tetrachloroethane	1	<
Bromomethane	2	<	2-Chloroethanol	5	<	1,2,3-Trichloropropane	1	<
Trichlorofluoromethane	1	<	1,2-Dichloropropane	1	<	Isopropylbenzene	1	<
Diethyl ether	5	<	Trichloroethene	1	<	Bromobenzene	1	<
Acetonitrile	5	<	Bromodichloromethane	1	<	2-Chlorotoluene	2	<
Idomethane	5	<	Dibromomethane	1	<	4-Chlorotoluene	2	<
1,1-Dichloroethene	1	<	Methyl methacrylate	5	<	n-propylbenzene	1	<
Carbon Disulfide	2	<	1,4-Dioxane	25	<	1,3,5-Trimethylbenzene	2	<
Allyl chloride	5	<	cis-1,3-Dichloropropene	1	<	Pentachloroethane	1	<
Methylene chloride	10	<	trans-1,3-Dichloropropene	1	<	1,2,4-Trimethylbenzene	2	<
Acrylonitrile	5	<	2-Nitropropane	5	<	tert-butylbenzene	1	<
1,1,2 - Trichlorotrifluoroethane	1	<	Toluene	1	<	sec-butylbenzene	1	<
trans-1,2-Dichloroethene	1	<	1,1,2-Trichloroethane	1	<	1,3-Dichlorobenzene	2	<
MTBE	1	<	Ethylmethacrylate	5	<	1,4-Dichlorobenzene	2	<
Chloroform	1	<	1,3-Dichloropropane	1	<	p-iso-Propyltoluene	1	<
1,1-Dichloroethane	1	<	Dibromochloromethane	1	<	1,2-Dichlorobenzene	2	<
Chloroprene	2	<	1,2-Dibromomethane	1	<	n-butylbenzene	1	<
2,2-Dichloropropane	1	<	Tetrachloroethene	1	<	1,2-Bromo-3-chloropropane	1	<
Methyl acrylate	5	<	Chlorobenzene	1	<	Nitrobenzene	5	<
cis-1,2,-Dichloroethene	1	<	1,1,1,2-Tetrachloroethane	1	<	1,2,4-Trichlorobenzene	1	<
Tetrahydrofuran	5	<	Ethylbenzene	1	<	Naphthalene	25	<
Methyacrylonitrile	5	<	m/p-xylene	1	<	1,2,3-Trichlorobenzene	1	<
Isobutyl alcohol	5	<	Bromoform	2	<	Hexachlorobutadiene	1	<
1,1,1-Trichloroethane	1	<	o-xylene	1	<			

Legend:
< = less than reporting limits
J = less than reporting limit, above detection levels
E = estimated, above calibration range

Surrogate Recovery	%
Dibromofluoromethane	95
Toluene-d8	104
4-Bromofluorobenzene	93

Notes:

Chemist/Reviewer: RMB

Date: 5/31/06

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Certificate of Analysis

Sample Identification	SB-27
Depth (ft)	0-4'
Sample Matrix	soil
Units:	ug/kg
% Solid	92.5%
Dilution	1
Sample Preservation	4 deg C.
Sample Volume	2 oz.
Date Sampled	5/25/2006
Date Analyzed	5/25/2006

Client:	**Woodard & Curran, Inc.**
Site:	**GP Mill**
Loc.:	**Old Town, Maine**
Project No.:	
Chemist:	D. M. Twomey

Analytical Method: USEPA SW-846 8260B
Instrument: Agilent 5890 GCD MS
Detectors: Mass spectrometer
Column: DB-5, 30 meter, .25 um
Conc. System: Tekmar 3000 Purge & Trap System
Trap: Vocarb 3000

Method:
USEPA SW846 METHOD 8260B

Compound	RL ug/kg	Results	Compound	RL ug/kg	Results	Compound	RL ug/kg	Results
Chloromethane	50	<	1,2-Dichloroethane	50	<	Styrene	50	<
Dichlorodifluoromethane	50	<	1,1-Dichloropropene	50	<	trans-1,4-Dichloro-2-butene	50	<
Chloroethane	50	<	Carbon Tetrachloride	50	<	cis-1,4,-Dichloro-2-butene	50	<
Vinyl chloride	50	<	Benzene	50	<	1,1,2,2-Tetrachloroethane	50	<
Bromomethane	100	<	2-Chloroethanol	250	<	1,2,3-Trichloropropane	50	<
Trichlorofluoromethane	50	<	1,2-Dichloropropane	50	<	Isopropylbenzene	50	<
Diethyl ether	250	<	Trichloroethene	50	<	Bromobenzene	50	<
Acetonitrile	250	<	Bromodichloromethane	50	<	2-Chlorotoluene	100	<
Idomethane	250	<	Dibromomethane	50	<	4-Chlorotoluene	100	<
1,1-Dichloroethene	50	<	Methyl methacrylate	250	<	n-propylbenzene	50	<
Carbon Disulfide	100	<	1,4-Dioxane	25	<	1,3,5-Trimethylbenzene	100	<
Allyl chloride	250	<	cis-1,3-Dichloropropene	50	<	Pentachloroethane	50	<
Methylene chloride	500	<	trans-1,3-Dichloropropene	50	<	1,2,4-Trimethylbenzene	100	<
Acrylonitrile	250	<	2-Nitropropane	250	<	tert-butylbenzene	50	<
1,1,2 - Trichlorotrifluoroethane	50	<	Toluene	50	<	sec-butylbenzene	50	<
trans-1,2-Dichloroethene	50	<	1,1,2-Trichloroethane	50	<	1,3-Dichlorobenzene	100	<
MTBE	50	<	Ethylmethacrylate	250	<	1,4-Dichlorobenzene	100	<
Chloroform	50	<	1,3-Dichloropropane	50	<	p-iso-Propyltoluene	50	<
1,1-Dichloroethane	50	<	Dibromochloromethane	50	<	1,2-Dichlorobenzene	100	<
Chloroprene	100	<	1,2-Dibromomethane	50	<	n-butylbenzene	50	<
2,2-Dichloropropane	50	<	Tetrachloroethene	50	<	1,2-Bromo-3-chloropropane	50	<
Methyl acrylate	250	<	Chlorobenzene	50	<	Nitrobenzene	250	<
cis-1,2,-Dichloroethene	50	<	1,1,1,2-Tetrachloroethane	50	<	1,2,4-Trichlorobenzene	50	<
Tetrahydrofuran	250	<	Ethylbenzene	50	<	Naphthalene	25	<
Methyacrylonitrile	250	<	m/p-xylene	50	<	1,2,3-Trichlorobenzene	50	<
Isobutyl alcohol	250	<	Bromoform	100	<	Hexachlorobutadiene	50	<
1,1,1-Trichloroethane	50	<	o-xylene	50	<			

Legend:
< = less than reporting limits
J = less than reporting limit, above detection levels
E = estimated, above calibration range

Surrogate Recovery	%
Dibromofluoromethane	77
Toluene-d8	111
4-Bromofluorobenzene	85

Notes:

Chemist/Reviewer: RMB

Date: 5/31/06

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Certificate of Analysis

Sample Identification	SB-29
Depth (ft)	0-4'
Sample Matrix	soil
Units:	ug/kg
% Solid	87.7%
Dilution	1
Sample Preservation	4 deg C.
Sample Volume	2 oz.
Date Sampled	5/25/2006
Date Analyzed	5/25/2006

Client:	**Woodard & Curran, Inc.**
Site:	**GP Mill**
Loc.:	**Old Town, Maine**
Project No.:	
Chemist:	D. M. Twomey

Analytical Method: USEPA SW-846 8260B
Instrument: Agilent 5890 GCD MS
Detectors: Mass spectrometer
Column: DB-5, 30 meter, .25 um
Conc. System: Tekmar 3000 Purge & Trap System
Trap: Vocarb 3000

Method:
USEPA SW846 METHOD 8260B

Compound	RL ug/kg	Results	Compound	RL ug/kg	Results	Compound	RL ug/kg	Results
Chloromethane	50	<	1,2-Dichloroethane	50	<	Styrene	50	<
Dichlorodifluoromethane	50	<	1,1-Dichloropropene	50	<	trans-1,4-Dichloro-2-butene	50	<
Chloroethane	50	<	Carbon Tetrachloride	50	<	cis-1,4,-Dichloro-2-butene	50	<
Vinyl chloride	50	<	Benzene	50	<	1,1,2,2-Tetrachloroethane	50	<
Bromomethane	100	<	2-Chloroethanol	250	<	1,2,3-Trichloropropane	50	<
Trichlorofluoromethane	50	<	1,2-Dichloropropane	50	<	Isopropylbenzene	50	<
Diethyl ether	250	<	Trichloroethene	50	<	Bromobenzene	50	<
Acetonitrile	250	<	Bromodichloromethane	50	<	2-Chlorotoluene	100	<
Idomethane	250	<	Dibromomethane	50	<	4-Chlorotoluene	100	<
1,1-Dichloroethene	50	<	Methyl methacrylate	250	<	n-propylbenzene	50	<
Carbon Disulfide	100	<	1,4-Dioxane	25	<	1,3,5-Trimethylbenzene	100	<
Allyl chloride	250	<	cis-1,3-Dichloropropene	50	<	Pentachloroethane	50	<
Methylene chloride	500	<	trans-1,3-Dichloropropene	50	<	1,2,4-Trimethylbenzene	100	<
Acrylonitrile	250	<	2-Nitropropane	250	<	tert-butylbenzene	50	<
1,1,2 - Trichlorotrifluoroethane	50	<	Toluene	50	<	sec-butylbenzene	50	<
trans-1,2-Dichloroethene	50	<	1,1,2-Trichloroethane	50	<	1,3-Dichlorobenzene	100	<
MTBE	50	<	Ethylmethacrylate	250	<	1,4-Dichlorobenzene	100	<
Chloroform	50	<	1,3-Dichloropropane	50	<	p-iso-Propyltoluene	50	<
1,1-Dichloroethane	50	<	Dibromochloromethane	50	<	1,2-Dichlorobenzene	100	<
Chloroprene	100	<	1,2-Dibromomethane	50	<	n-butylbenzene	50	<
2,2-Dichloropropane	50	<	Tetrachloroethene	50	<	1,2-Bromo-3-chloropropane	50	<
Methyl acrylate	250	<	Chlorobenzene	50	<	Nitrobenzene	250	<
cis-1,2,-Dichloroethene	50	<	1,1,1,2-Tetrachloroethane	50	<	1,2,4-Trichlorobenzene	50	<
Tetrahydrofuran	250	<	Ethylbenzene	50	<	Naphthalene	25	<
Methyacrylonitrile	250	<	m/p-xylene	50	<	1,2,3-Trichlorobenzene	50	<
Isobutyl alcohol	250	<	Bromoform	100	<	Hexachlorobutadiene	50	<
1,1,1-Trichloroethane	50	<	o-xylene	50	<			

Legend:
< = less than reporting limits
J = less than reporting limit, above detection levels
E = estimated, above calibration range

Surrogate Recovery	%
Dibromofluoromethane	110
Toluene-d8	104
4-Bromofluorobenzene	82

Notes:

Chemist/Reviewer: RMB

Date: 5/31/06

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Certificate of Analysis

Sample Identification	SB-32
Depth (ft)	6-10'
Sample Matrix	water
Units:	ug/L
% Solid	NA
Dilution	1
Sample Preservation	4 deg C.
Sample Volume	40 mL.
Date Sampled	5/25/2006
Date Analyzed	5/25/2006

Client:	**Woodard & Curran, Inc.**
Site:	**GP Mill**
Loc.:	**Old Town, Maine**
Project No.:	
Chemist:	D. M. Twomey

Analytical Method: USEPA SW-846 8260B
Instrument: Agilent 5890 GCD MS
Detectors: Mass spectrometer
Column: DB-5, 30 meter, .25 um
Conc. System: Tekmar 3000 Purge & Trap System
Trap: Vocarb 3000

Method:
USEPA SW846 METHOD 8260B

Compound	RL ug/L	Results	Compound	RL ug/L	Results	Compound	RL ug/L	Results
Chloromethane	1	<	1,2-Dichloroethane	1	<	Styrene	1	<
Dichlorodifluoromethane	1	<	1,1-Dichloropropene	1	<	trans-1,4-Dichloro-2-butene	1	<
Chloroethane	1	<	Carbon Tetrachloride	1	<	cis-1,4,-Dichloro-2-butene	1	<
Vinyl chloride	1	<	Benzene	1	2.1	1,1,2,2-Tetrachloroethane	1	<
Bromomethane	2	<	2-Chloroethanol	5	<	1,2,3-Trichloropropane	1	<
Trichlorofluoromethane	1	<	1,2-Dichloropropane	1	<	Isopropylbenzene	1	<
Diethyl ether	5	<	Trichloroethene	1	<	Bromobenzene	1	<
Acetonitrile	5	<	Bromodichloromethane	1	<	2-Chlorotoluene	2	<
Idomethane	5	<	Dibromomethane	1	<	4-Chlorotoluene	2	<
1,1-Dichloroethene	1	<	Methyl methacrylate	5	<	n-propylbenzene	1	<
Carbon Disulfide	2	<	1,4-Dioxane	25	<	1,3,5-Trimethylbenzene	2	<
Allyl chloride	5	<	cis-1,3-Dichloropropene	1	<	Pentachloroethane	1	<
Methylene chloride	10	<	trans-1,3-Dichloropropene	1	<	1,2,4-Trimethylbenzene	2	2.9
Acrylonitrile	5	<	2-Nitropropane	5	<	tert-butylbenzene	1	<
1,1,2 - Trichlorotrifluoroethane	1	<	Toluene	1	<	sec-butylbenzene	1	<
trans-1,2-Dichloroethene	1	<	1,1,2-Trichloroethane	1	<	1,3-Dichlorobenzene	2	<
MTBE	1	<	Ethylmethacrylate	5	<	1,4-Dichlorobenzene	2	<
Chloroform	1	<	1,3-Dichloropropane	1	<	p-iso-Propyltoluene	1	<
1,1-Dichloroethane	1	<	Dibromochloromethane	1	<	1,2-Dichlorobenzene	2	<
Chloroprene	2	<	1,2-Dibromomethane	1	<	n-butylbenzene	1	<
2,2-Dichloropropane	1	<	Tetrachloroethene	1	<	1,2-Bromo-3-chloropropane	1	<
Methyl acrylate	5	<	Chlorobenzene	1	<	Nitrobenzene	5	<
cis-1,2,-Dichloroethene	1	<	1,1,1,2-Tetrachloroethane	1	<	1,2,4-Trichlorobenzene	1	<
Tetrahydrofuran	5	<	Ethylbenzene	1	<	Naphthalene	25	<
Methyacrylonitrile	5	<	m/p-xylene	1	<	1,2,3-Trichlorobenzene	1	<
Isobutyl alcohol	5	<	Bromoform	2	<	Hexachlorobutadiene	1	<
1,1,1-Trichloroethane	1	<	o-xylene	1	<			

Legend:
< = less than reporting limits
J = less than reporting limit, above detection levels
E = estimated, above calibration range

Surrogate Recovery	%
Dibromofluoromethane	91
Toluene-d8	107
4-Bromofluorobenzene	95

Notes:

Chemist/Reviewer: RMB

Date: 5/31/06

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Sample Identification	SB-33
Depth (ft)	6-10'
Sample Matrix	water
Units:	ug/L
% Solid	NA
Dilution	1
Sample Preservation	4 deg C.
Sample Volume	40 mL.
Date Sampled	5/25/2006
Date Analyzed	5/25/2006

Client:	**Woodard & Curran, Inc.**
Site:	**GP Mill**
Loc.:	**Old Town, Maine**
Project No.:	
Chemist:	D. M. Twomey

Analytical Method: USEPA SW-846 8260B
Instrument: Agilent 5890 GCD MS
Detectors: Mass spectrometer
Column: DB-5, 30 meter, .25 um
Conc. System: Tekmar 3000 Purge & Trap System
Trap: Vocarb 3000

Method:
USEPA SW846 METHOD 8260B

Compound	RL ug/L	Results	Compound	RL ug/L	Results	Compound	RL ug/L	Results
Chloromethane	1	<	1,2-Dichloroethane	1	<	Styrene	1	<
Dichlorodifluoromethane	1	<	1,1-Dichloropropene	1	<	trans-1,4-Dichloro-2-butene	1	<
Chloroethane	1	<	Carbon Tetrachloride	1	<	cis-1,4,-Dichloro-2-butene	1	<
Vinyl chloride	1	<	Benzene	1	<	1,1,2,2-Tetrachloroethane	1	<
Bromomethane	2	<	2-Chloroethanol	5	<	1,2,3-Trichloropropane	1	<
Trichlorofluoromethane	1	<	1,2-Dichloropropane	1	<	Isopropylbenzene	1	<
Diethyl ether	5	<	Trichloroethene	1	<	Bromobenzene	1	<
Acetonitrile	5	<	Bromodichloromethane	1	<	2-Chlorotoluene	2	<
Idomethane	5	<	Dibromomethane	1	<	4-Chlorotoluene	2	<
1,1-Dichloroethene	1	<	Methyl methacrylate	5	<	n-propylbenzene	1	<
Carbon Disulfide	2	<	1,4-Dioxane	25	<	1,3,5-Trimethylbenzene	2	<
Allyl chloride	5	<	cis-1,3-Dichloropropene	1	<	Pentachloroethane	1	<
Methylene chloride	10	<	trans-1,3-Dichloropropene	1	<	1,2,4-Trimethylbenzene	2	<
Acrylonitrile	5	<	2-Nitropropane	5	<	tert-butylbenzene	1	<
1,1,2 - Trichlorotrifluoroethane	1	<	Toluene	1	<	sec-butylbenzene	1	<
trans-1,2-Dichloroethene	1	<	1,1,2-Trichloroethane	1	<	1,3-Dichlorobenzene	2	<
MTBE	1	<	Ethylmethacrylate	5	<	1,4-Dichlorobenzene	2	<
Chloroform	1	<	1,3-Dichloropropane	1	<	p-iso-Propyltoluene	1	<
1,1-Dichloroethane	1	<	Dibromochloromethane	1	<	1,2-Dichlorobenzene	2	<
Chloroprene	2	<	1,2-Dibromomethane	1	<	n-butylbenzene	1	<
2,2-Dichloropropane	1	<	Tetrachloroethene	1	<	1,2-Bromo-3-chloropropane	1	<
Methyl acrylate	5	<	Chlorobenzene	1	<	Nitrobenzene	5	<
cis-1,2,-Dichloroethene	1	<	1,1,1,2-Tetrachloroethane	1	<	1,2,4-Trichlorobenzene	1	<
Tetrahydrofuran	5	<	Ethylbenzene	1	<	Naphthalene	25	<
Methyacrylonitrile	5	<	m/p-xylene	1	<	1,2,3-Trichlorobenzene	1	<
Isobutyl alcohol	5	<	Bromoform	2	<	Hexachlorobutadiene	1	<
1,1,1-Trichloroethane	1	<	o-xylene	1	<	*Chlorodifluorethane***	*1*	**13**

Legend:
< = less than reporting limits
J = less than reporting limit, above detection levels
E = estimated, above calibration range

Surrogate Recovery	%
Dibromofluoromethane	94
Toluene-d8	103
4-Bromofluorobenzene	92

Notes:
** - **non-target compound identified.**
Sample had non-target chromatography characteristics of No. 2/diesel fuel present.

Chemist/Reviewer: RMB

Date: 5/31/06

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Sample Identification	SB-15
Depth (ft)	3-8'
Sample Matrix	water
Units:	ug/L
% Solid	NA
Dilution	1
Sample Preservation	4 deg C.
Sample Volume	40 mL.
Date Sampled	5/25/2006
Date Analyzed	5/25/2006

Client:	**Woodard & Curran, Inc.**
Site:	**GP Mill**
Loc.:	**Old Town, Maine**
Project No.:	
Chemist:	D. M. Twomey

Analytical Method: USEPA SW-846 8260B
Instrument: Agilent 5890 GCD MS
Detectors: Mass spectrometer
Column: DB-5, 30 meter, .25 um
Conc. System: Tekmar 3000 Purge & Trap System
Trap: Vocarb 3000

Method:
USEPA SW846 METHOD 8260B

Compound	RL ug/L	Results	Compound	RL ug/L	Results	Compound	RL ug/L	Results
Chloromethane	1	<	1,2-Dichloroethane	1	<	Styrene	1	<
Dichlorodifluoromethane	1	<	1,1-Dichloropropene	1	<	trans-1,4-Dichloro-2-butene	1	<
Chloroethane	1	<	Carbon Tetrachloride	1	<	cis-1,4,-Dichloro-2-butene	1	<
Vinyl chloride	1	<	Benzene	1	<	1,1,2,2-Tetrachloroethane	1	<
Bromomethane	2	<	2-Chloroethanol	5	<	1,2,3-Trichloropropane	1	<
Trichlorofluoromethane	1	<	1,2-Dichloropropane	1	<	Isopropylbenzene	1	<
Diethyl ether	5	<	Trichloroethene	1	<	Bromobenzene	1	<
Acetonitrile	5	<	Bromodichloromethane	1	<	2-Chlorotoluene	2	<
Idomethane	5	<	Dibromomethane	1	<	4-Chlorotoluene	2	<
1,1-Dichloroethene	1	<	Methyl methacrylate	5	<	n-propylbenzene	1	<
Carbon Disulfide	2	<	1,4-Dioxane	25	<	1,3,5-Trimethylbenzene	2	<
Allyl chloride	5	<	cis-1,3-Dichloropropene	1	<	Pentachloroethane	1	<
Methylene chloride	10	<	trans-1,3-Dichloropropene	1	<	1,2,4-Trimethylbenzene	2	<
Acrylonitrile	5	<	2-Nitropropane	5	<	tert-butylbenzene	1	<
1,1,2 - Trichlorotrifluoroethane	1	<	Toluene	1	<	sec-butylbenzene	1	<
trans-1,2-Dichloroethene	1	<	1,1,2-Trichloroethane	1	<	1,3-Dichlorobenzene	2	<
MTBE	1	<	Ethylmethacrylate	5	<	1,4-Dichlorobenzene	2	<
Chloroform	1	<	1,3-Dichloropropane	1	<	p-iso-Propyltoluene	1	<
1,1-Dichloroethane	1	<	Dibromochloromethane	1	<	1,2-Dichlorobenzene	2	<
Chloroprene	2	<	1,2-Dibromomethane	1	<	n-butylbenzene	1	<
2,2-Dichloropropane	1	<	Tetrachloroethene	1	<	1,2-Bromo-3-chloropropane	1	<
Methyl acrylate	5	<	Chlorobenzene	1	<	Nitrobenzene	5	<
cis-1,2,-Dichloroethene	1	<	1,1,1,2-Tetrachloroethane	1	<	1,2,4-Trichlorobenzene	1	<
Tetrahydrofuran	5	<	Ethylbenzene	1	<	Naphthalene	25	<
Methyacrylonitrile	5	<	m/p-xylene	1	<	1,2,3-Trichlorobenzene	1	<
Isobutyl alcohol	5	<	Bromoform	2	<	Hexachlorobutadiene	1	<
1,1,1-Trichloroethane	1	<	o-xylene	1	<			

Legend:
< = less than reporting limits
J = less than reporting limit, above detection levels
E = estimated, above calibration range

Surrogate Recovery	%
Dibromofluoromethane	87
Toluene-d8	100
4-Bromofluorobenzene	97

Notes:

Chemist/Reviewer: RMB

Date: 5/31/06

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Certificate of Analysis

Sample Identification	SB-26
Depth (ft)	8-12'
Sample Matrix	water
Units:	ug/L
% Solid	NA
Dilution	1
Sample Preservation	4 deg C.
Sample Volume	40 mL.
Date Sampled	5/25/2006
Date Analyzed	5/25/2006

Client:	**Woodard & Curran, Inc.**
Site:	**GP Mill**
Loc.:	**Old Town, Maine**
Project No.:	
Chemist:	D. M. Twomey

Analytical Method: USEPA SW-846 8260B
Instrument: Agilent 5890 GCD MS
Detectors: Mass spectrometer
Column: DB-5, 30 meter, .25 um
Conc. System: Tekmar 3000 Purge & Trap System
Trap: Vocarb 3000

Method:
USEPA SW846 METHOD 8260B

Compound	RL ug/L	Results	Compound	RL ug/L	Results	Compound	RL ug/L	Results
Chloromethane	1	<	1,2-Dichloroethane	1	<	Styrene	1	<
Dichlorodifluoromethane	1	<	1,1-Dichloropropene	1	<	trans-1,4-Dichloro-2-butene	1	<
Chloroethane	1	<	Carbon Tetrachloride	1	<	cis-1,4,-Dichloro-2-butene	1	<
Vinyl chloride	1	<	Benzene	1	<	1,1,2,2-Tetrachloroethane	1	<
Bromomethane	2	<	2-Chloroethanol	5	<	1,2,3-Trichloropropane	1	<
Trichlorofluoromethane	1	<	1,2-Dichloropropane	1	<	Isopropylbenzene	1	<
Diethyl ether	5	<	Trichloroethene	1	<	Bromobenzene	1	<
Acetonitrile	5	<	Bromodichloromethane	1	<	2-Chlorotoluene	2	<
Idomethane	5	<	Dibromomethane	1	<	4-Chlorotoluene	2	<
1,1-Dichloroethene	1	<	Methyl methacrylate	5	<	n-propylbenzene	1	<
Carbon Disulfide	2	<	1,4-Dioxane	25	<	1,3,5-Trimethylbenzene	2	<
Allyl chloride	5	<	cis-1,3-Dichloropropene	1	<	Pentachloroethane	1	<
Methylene chloride	10	<	trans-1,3-Dichloropropene	1	<	1,2,4-Trimethylbenzene	2	<
Acrylonitrile	5	<	2-Nitropropane	5	<	tert-butylbenzene	1	<
1,1,2 - Trichlorotrifluoroethane	1	<	Toluene	1	<	sec-butylbenzene	1	<
trans-1,2-Dichloroethene	1	<	1,1,2-Trichloroethane	1	<	1,3-Dichlorobenzene	2	<
MTBE	1	<	Ethylmethacrylate	5	<	1,4-Dichlorobenzene	2	<
Chloroform	1	<	1,3-Dichloropropane	1	<	p-iso-Propyltoluene	1	<
1,1-Dichloroethane	1	<	Dibromochloromethane	1	<	1,2-Dichlorobenzene	2	<
Chloroprene	2	<	1,2-Dibromomethane	1	<	n-butylbenzene	1	<
2,2-Dichloropropane	1	<	Tetrachloroethene	1	<	1,2-Bromo-3-chloropropane	1	<
Methyl acrylate	5	<	Chlorobenzene	1	<	Nitrobenzene	5	<
cis-1,2,-Dichloroethene	1	<	1,1,1,2-Tetrachloroethane	1	<	1,2,4-Trichlorobenzene	1	<
Tetrahydrofuran	5	<	Ethylbenzene	1	<	Naphthalene	25	<
Methyacrylonitrile	5	<	m/p-xylene	1	<	1,2,3-Trichlorobenzene	1	<
Isobutyl alcohol	5	<	Bromoform	2	<	Hexachlorobutadiene	1	<
1,1,1-Trichloroethane	1	<	o-xylene	1	<			

Legend:
< = less than reporting limits
J = less than reporting limit, above detection levels
E = estimated, above calibration range

Surrogate Recovery	%
Dibromofluoromethane	88
Toluene-d8	100
4-Bromofluorobenzene	98

Notes:

Chemist/Reviewer: RMB

Date: 5/31/06

PO Box 6752
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15 Holly St., Ste 102
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800-880-0533
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Certificate of Analysis

Sample Identification	SB-25
Depth (ft)	7-11'
Sample Matrix	water
Units:	ug/L
% Solid	NA
Dilution	1
Sample Preservation	4 deg C.
Sample Volume	40 mL.
Date Sampled	5/25/2006
Date Analyzed	5/25/2006

Client:	**Woodard & Curran, Inc.**
Site:	**GP Mill**
Loc.:	**Old Town, Maine**
Project No.:	
Chemist:	D. M. Twomey

Analytical Method: USEPA SW-846 8260B
Instrument: Agilent 5890 GCD MS
Detectors: Mass spectrometer
Column: DB-5, 30 meter, .25 um
Conc. System: Tekmar 3000 Purge & Trap System
Trap: Vocarb 3000

Method:
USEPA SW846 METHOD 8260B

Compound	RL ug/L	Results	Compound	RL ug/L	Results	Compound	RL ug/L	Results
Chloromethane	1	<	1,2-Dichloroethane	1	<	Styrene	1	<
Dichlorodifluoromethane	1	<	1,1-Dichloropropene	1	<	trans-1,4-Dichloro-2-butene	1	<
Chloroethane	1	<	Carbon Tetrachloride	1	<	cis-1,4,-Dichloro-2-butene	1	<
Vinyl chloride	1	<	Benzene	1	<	1,1,2,2-Tetrachloroethane	1	<
Bromomethane	2	<	2-Chloroethanol	5	<	1,2,3-Trichloropropane	1	<
Trichlorofluoromethane	1	<	1,2-Dichloropropane	1	<	Isopropylbenzene	1	<
Diethyl ether	5	<	Trichloroethene	1	<	Bromobenzene	1	<
Acetonitrile	5	<	Bromodichloromethane	1	<	2-Chlorotoluene	2	<
Idomethane	5	<	Dibromomethane	1	<	4-Chlorotoluene	2	<
1,1-Dichloroethene	1	<	Methyl methacrylate	5	<	n-propylbenzene	1	<
Carbon Disulfide	2	<	1,4-Dioxane	25	<	1,3,5-Trimethylbenzene	2	<
Allyl chloride	5	<	cis-1,3-Dichloropropene	1	<	Pentachloroethane	1	<
Methylene chloride	10	<	trans-1,3-Dichloropropene	1	<	1,2,4-Trimethylbenzene	2	<
Acrylonitrile	5	<	2-Nitropropane	5	<	tert-butylbenzene	1	<
1,1,2 - Trichlorotrifluoroethane	1	<	Toluene	1	<	sec-butylbenzene	1	<
trans-1,2-Dichloroethene	1	<	1,1,2-Trichloroethane	1	<	1,3-Dichlorobenzene	2	<
MTBE	1	<	Ethylmethacrylate	5	<	1,4-Dichlorobenzene	2	<
Chloroform	1	<	1,3-Dichloropropane	1	<	p-iso-Propyltoluene	1	<
1,1-Dichloroethane	1	<	Dibromochloromethane	1	<	1,2-Dichlorobenzene	2	<
Chloroprene	2	<	1,2-Dibromomethane	1	<	n-butylbenzene	1	<
2,2-Dichloropropane	1	<	Tetrachloroethene	1	<	1,2-Bromo-3-chloropropane	1	<
Methyl acrylate	5	<	Chlorobenzene	1	<	Nitrobenzene	5	<
cis-1,2,-Dichloroethene	1	<	1,1,1,2-Tetrachloroethane	1	<	1,2,4-Trichlorobenzene	1	<
Tetrahydrofuran	5	<	Ethylbenzene	1	<	Naphthalene	25	<
Methyacrylonitrile	5	<	m/p-xylene	1	<	1,2,3-Trichlorobenzene	1	<
Isobutyl alcohol	5	<	Bromoform	2	<	Hexachlorobutadiene	1	<
1,1,1-Trichloroethane	1	<	o-xylene	1	<			

Legend:
< = less than reporting limits
J = less than reporting limit, above detection levels
E = estimated, above calibration range

Surrogate Recovery	%
Dibromofluoromethane	87
Toluene-d8	107
4-Bromofluorobenzene	92

Notes:

Chemist/Reviewer: RMB

Date: 5/31/06

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Certificate of Analysis

Sample Identification	SB-30
Depth (ft)	6-10'
Sample Matrix	water
Units:	ug/L
% Solid	NA
Dilution	1
Sample Preservation	4 deg C.
Sample Volume	40 mL.
Date Sampled	5/25/2006
Date Analyzed	5/25/2006

Client:	**Woodard & Curran, Inc.**
Site:	**GP Mill**
Loc.:	**Old Town, Maine**
Project No.:	
Chemist:	D. M. Twomey

Analytical Method: USEPA SW-846 8260B
Instrument: Agilent 5890 GCD MS
Detectors: Mass spectrometer
Column: DB-5, 30 meter, .25 um
Conc. System: Tekmar 3000 Purge & Trap System
Trap: Vocarb 3000

Method:
USEPA SW846 METHOD 8260B

Compound	RL ug/L	Results	Compound	RL ug/L	Results	Compound	RL ug/L	Results
Chloromethane	1	<	1,2-Dichloroethane	1	<	Styrene	1	<
Dichlorodifluoromethane	1	<	1,1-Dichloropropene	1	<	trans-1,4-Dichloro-2-butene	1	<
Chloroethane	1	<	Carbon Tetrachloride	1	<	cis-1,4,-Dichloro-2-butene	1	<
Vinyl chloride	1	<	Benzene	1	<	1,1,2,2-Tetrachloroethane	1	<
Bromomethane	2	<	2-Chloroethanol	5	<	1,2,3-Trichloropropane	1	<
Trichlorofluoromethane	1	<	1,2-Dichloropropane	1	<	Isopropylbenzene	1	<
Diethyl ether	5	<	Trichloroethene	1	<	Bromobenzene	1	<
Acetonitrile	5	<	Bromodichloromethane	1	<	2-Chlorotoluene	2	<
Idomethane	5	<	Dibromomethane	1	<	4-Chlorotoluene	2	<
1,1-Dichloroethene	1	<	Methyl methacrylate	5	<	n-propylbenzene	1	<
Carbon Disulfide	2	<	1,4-Dioxane	25	<	1,3,5-Trimethylbenzene	2	<
Allyl chloride	5	<	cis-1,3-Dichloropropene	1	<	Pentachloroethane	1	<
Methylene chloride	10	<	trans-1,3-Dichloropropene	1	<	1,2,4-Trimethylbenzene	2	<
Acrylonitrile	5	<	2-Nitropropane	5	<	tert-butylbenzene	1	<
1,1,2 - Trichlorotrifluoroethane	1	<	Toluene	1	<	sec-butylbenzene	1	<
trans-1,2-Dichloroethene	1	<	1,1,2-Trichloroethane	1	<	1,3-Dichlorobenzene	2	<
MTBE	1	<	Ethylmethacrylate	5	<	1,4-Dichlorobenzene	2	<
Chloroform	1	<	1,3-Dichloropropane	1	<	p-iso-Propyltoluene	1	<
1,1-Dichloroethane	1	<	Dibromochloromethane	1	<	1,2-Dichlorobenzene	2	<
Chloroprene	2	<	1,2-Dibromomethane	1	<	n-butylbenzene	1	<
2,2-Dichloropropane	1	<	Tetrachloroethene	1	<	1,2-Bromo-3-chloropropane	1	<
Methyl acrylate	5	<	Chlorobenzene	1	<	Nitrobenzene	5	<
cis-1,2,-Dichloroethene	1	<	1,1,1,2-Tetrachloroethane	1	<	1,2,4-Trichlorobenzene	1	<
Tetrahydrofuran	5	<	Ethylbenzene	1	<	Naphthalene	25	<
Methyacrylonitrile	5	<	m/p-xylene	1	<	1,2,3-Trichlorobenzene	1	<
Isobutyl alcohol	5	<	Bromoform	2	<	Hexachlorobutadiene	1	<
1,1,1-Trichloroethane	1	<	o-xylene	1	<			

Legend:
< = less than reporting limits
J = less than reporting limit, above detection levels
E = estimated, above calibration range

Surrogate Recovery	%
Dibromofluoromethane	90
Toluene-d8	107
4-Bromofluorobenzene	93

Notes:

Chemist/Reviewer: RMB

Date: 5/31/06

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Certificate of Analysis

Sample Identification	SB-31
Depth (ft)	4-8'
Sample Matrix	water
Units:	ug/L
% Solid	NA
Dilution	1
Sample Preservation	4 deg C.
Sample Volume	40 mL.
Date Sampled	5/25/2006
Date Analyzed	5/25/2006

Client:	Woodard & Curran, Inc.
Site:	GP Mill
Loc.:	Old Town, Maine
Project No.:	
Chemist:	D. M. Twomey

Analytical Method: USEPA SW-846 8260B
Instrument: Agilent 5890 GCD MS
Detectors: Mass spectrometer
Column: DB-5, 30 meter, .25 um
Conc. System: Tekmar 3000 Purge & Trap System
Trap: Vocarb 3000

Method:
USEPA SW846 METHOD 8260B

Compound	RL ug/L	Results	Compound	RL ug/L	Results	Compound	RL ug/L	Results
Chloromethane	1	<	1,2-Dichloroethane	1	<	Styrene	1	<
Dichlorodifluoromethane	1	<	1,1-Dichloropropene	1	<	trans-1,4-Dichloro-2-butene	1	<
Chloroethane	1	<	Carbon Tetrachloride	1	<	cis-1,4,-Dichloro-2-butene	1	<
Vinyl chloride	1	<	Benzene	1	<	1,1,2,2-Tetrachloroethane	1	<
Bromomethane	2	<	2-Chloroethanol	5	<	1,2,3-Trichloropropane	1	<
Trichlorofluoromethane	1	<	1,2-Dichloropropane	1	<	Isopropylbenzene	1	<
Diethyl ether	5	<	Trichloroethene	1	<	Bromobenzene	1	<
Acetonitrile	5	<	Bromodichloromethane	1	<	2-Chlorotoluene	2	<
Idomethane	5	<	Dibromomethane	1	<	4-Chlorotoluene	2	<
1,1-Dichloroethene	1	<	Methyl methacrylate	5	<	n-propylbenzene	1	<
Carbon Disulfide	2	<	1,4-Dioxane	25	<	1,3,5-Trimethylbenzene	2	<
Allyl chloride	5	<	cis-1,3-Dichloropropene	1	<	Pentachloroethane	1	<
Methylene chloride	10	<	trans-1,3-Dichloropropene	1	<	1,2,4-Trimethylbenzene	2	<
Acrylonitrile	5	<	2-Nitropropane	5	<	tert-butylbenzene	1	<
1,1,2 - Trichlorotrifluoroethane	1	<	Toluene	1	<	sec-butylbenzene	1	<
trans-1,2-Dichloroethene	1	<	1,1,2-Trichloroethane	1	<	1,3-Dichlorobenzene	2	<
MTBE	1	<	Ethylmethacrylate	5	<	1,4-Dichlorobenzene	2	<
Chloroform	1	<	1,3-Dichloropropane	1	<	p-iso-Propyltoluene	1	<
1,1-Dichloroethane	1	<	Dibromochloromethane	1	<	1,2-Dichlorobenzene	2	<
Chloroprene	2	<	1,2-Dibromomethane	1	<	n-butylbenzene	1	<
2,2-Dichloropropane	1	<	Tetrachloroethene	1	<	1,2-Bromo-3-chloropropane	1	<
Methyl acrylate	5	<	Chlorobenzene	1	<	Nitrobenzene	5	<
cis-1,2,-Dichloroethene	1	<	1,1,1,2-Tetrachloroethane	1	<	1,2,4-Trichlorobenzene	1	<
Tetrahydrofuran	5	<	Ethylbenzene	1	<	Naphthalene	25	<
Methyacrylonitrile	5	<	m/p-xylene	1	<	1,2,3-Trichlorobenzene	1	<
Isobutyl alcohol	5	<	Bromoform	2	<	Hexachlorobutadiene	1	<
1,1,1-Trichloroethane	1	<	o-xylene	1	<			

Legend:
< = less than reporting limits
J = less than reporting limit, above detection levels
E = estimated, above calibration range

Surrogate Recovery	%
Dibromofluoromethane	91
Toluene-d8	101
4-Bromofluorobenzene	96

Notes:

Chemist/Reviewer: RMB

Date: 5/31/06

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Certificate of Analysis

Sample Identification	SB-38
Depth (ft)	8-12'
Sample Matrix	water
Units:	ug/L
% Solid	NA
Dilution	1
Sample Preservation	4 deg C.
Sample Volume	40 mL.
Date Sampled	5/25/2006
Date Analyzed	5/25/2006

Client:	**Woodard & Curran, Inc.**
Site:	**GP Mill**
Loc.:	**Old Town, Maine**
Project No.:	
Chemist:	D. M. Twomey

Analytical Method: USEPA SW-846 8260B
Instrument: Agilent 5890 GCD MS
Detectors: Mass spectrometer
Column: DB-5, 30 meter, .25 um
Conc. System: Tekmar 3000 Purge & Trap System
Trap: Vocarb 3000

Method:
USEPA SW846 METHOD 8260B

Compound	RL ug/L	Results	Compound	RL ug/L	Results	Compound	RL ug/L	Results
Chloromethane	1	<	1,2-Dichloroethane	1	<	Styrene	1	<
Dichlorodifluoromethane	1	<	1,1-Dichloropropene	1	<	trans-1,4-Dichloro-2-butene	1	<
Chloroethane	1	<	Carbon Tetrachloride	1	<	cis-1,4,-Dichloro-2-butene	1	<
Vinyl chloride	1	<	Benzene	1	<	1,1,2,2-Tetrachloroethane	1	<
Bromomethane	2	<	2-Chloroethanol	5	<	1,2,3-Trichloropropane	1	<
Trichlorofluoromethane	1	<	1,2-Dichloropropane	1	<	Isopropylbenzene	1	<
Diethyl ether	5	<	Trichloroethene	1	<	Bromobenzene	1	<
Acetonitrile	5	<	Bromodichloromethane	1	<	2-Chlorotoluene	2	<
Idomethane	5	<	Dibromomethane	1	<	4-Chlorotoluene	2	<
1,1-Dichloroethene	1	<	Methyl methacrylate	5	<	n-propylbenzene	1	<
Carbon Disulfide	2	<	1,4-Dioxane	25	<	1,3,5-Trimethylbenzene	2	<
Allyl chloride	5	<	cis-1,3-Dichloropropene	1	<	Pentachloroethane	1	<
Methylene chloride	10	<	trans-1,3-Dichloropropene	1	<	1,2,4-Trimethylbenzene	2	<
Acrylonitrile	5	<	2-Nitropropane	5	<	tert-butylbenzene	1	<
1,1,2 - Trichlorotrifluoroethane	1	<	Toluene	1	<	sec-butylbenzene	1	<
trans-1,2-Dichloroethene	1	<	1,1,2-Trichloroethane	1	<	1,3-Dichlorobenzene	2	<
MTBE	1	<	Ethylmethacrylate	5	<	1,4-Dichlorobenzene	2	<
Chloroform	1	<	1,3-Dichloropropane	1	<	p-iso-Propyltoluene	1	<
1,1-Dichloroethane	1	<	Dibromochloromethane	1	<	1,2-Dichlorobenzene	2	<
Chloroprene	2	<	1,2-Dibromomethane	1	<	n-butylbenzene	1	<
2,2-Dichloropropane	1	<	Tetrachloroethene	1	<	1,2-Bromo-3-chloropropane	1	<
Methyl acrylate	5	<	Chlorobenzene	1	<	Nitrobenzene	5	<
cis-1,2,-Dichloroethene	1	<	1,1,1,2-Tetrachloroethane	1	<	1,2,4-Trichlorobenzene	1	<
Tetrahydrofuran	5	<	Ethylbenzene	1	<	Naphthalene	25	<
Methyacrylonitrile	5	<	m/p-xylene	1	<	1,2,3-Trichlorobenzene	1	<
Isobutyl alcohol	5	<	Bromoform	2	<	Hexachlorobutadiene	1	<
1,1,1-Trichloroethane	1	<	o-xylene	1	<			

Legend:
< = less than reporting limits
J = less than reporting limit, above detection levels
E = estimated, above calibration range

Surrogate Recovery	%
Dibromofluoromethane	78
Toluene-d8	101
4-Bromofluorobenzene	86

Notes:

Chemist/Reviewer: RMB

Date: 5/31/06

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Certificate of Analysis

Sample Identification	SB-37
Depth (ft)	7.3-11.7'
Sample Matrix	water
Units:	ug/L
% Solid	NA
Dilution	1
Sample Preservation	4 deg C.
Sample Volume	40 mL.
Date Sampled	5/25/2006
Date Analyzed	5/25/2006

Client:	Woodard & Curran, Inc.
Site:	GP Mill
Loc.:	Old Town, Maine
Project No.:	
Chemist:	D. M. Twomey

Analytical Method: USEPA SW-846 8260B
Instrument: Agilent 5890 GCD MS
Detectors: Mass spectrometer
Column: DB-5, 30 meter, .25 um
Conc. System: Tekmar 3000 Purge & Trap System
Trap: Vocarb 3000

Method:
USEPA SW846 METHOD 8260B

Compound	RL ug/L	Results	Compound	RL ug/L	Results	Compound	RL ug/L	Results
Chloromethane	1	<	1,2-Dichloroethane	1	<	Styrene	1	<
Dichlorodifluoromethane	1	<	1,1-Dichloropropene	1	<	trans-1,4-Dichloro-2-butene	1	<
Chloroethane	1	<	Carbon Tetrachloride	1	<	cis-1,4,-Dichloro-2-butene	1	<
Vinyl chloride	1	<	Benzene	1	<	1,1,2,2-Tetrachloroethane	1	<
Bromomethane	2	<	2-Chloroethanol	5	<	1,2,3-Trichloropropane	1	<
Trichlorofluoromethane	1	<	1,2-Dichloropropane	1	<	Isopropylbenzene	1	<
Diethyl ether	5	<	Trichloroethene	1	<	Bromobenzene	1	<
Acetonitrile	5	<	Bromodichloromethane	1	<	2-Chlorotoluene	2	<
Idomethane	5	<	Dibromomethane	1	<	4-Chlorotoluene	2	<
1,1-Dichloroethene	1	<	Methyl methacrylate	5	<	n-propylbenzene	1	<
Carbon Disulfide	2	<	1,4-Dioxane	25	<	1,3,5-Trimethylbenzene	2	<
Allyl chloride	5	<	cis-1,3-Dichloropropene	1	<	Pentachloroethane	1	<
Methylene chloride	10	<	trans-1,3-Dichloropropene	1	<	1,2,4-Trimethylbenzene	2	<
Acrylonitrile	5	<	2-Nitropropane	5	<	tert-butylbenzene	1	<
1,1,2 - Trichlorotrifluoroethane	1	<	Toluene	1	<	sec-butylbenzene	1	<
trans-1,2-Dichloroethene	1	<	1,1,2-Trichloroethane	1	<	1,3-Dichlorobenzene	2	<
MTBE	1	<	Ethylmethacrylate	5	<	1,4-Dichlorobenzene	2	<
Chloroform	1	<	1,3-Dichloropropane	1	<	p-iso-Propyltoluene	1	<
1,1-Dichloroethane	1	<	Dibromochloromethane	1	<	1,2-Dichlorobenzene	2	<
Chloroprene	2	<	1,2-Dibromomethane	1	<	n-butylbenzene	1	<
2,2-Dichloropropane	1	<	Tetrachloroethene	1	<	1,2-Bromo-3-chloropropane	1	<
Methyl acrylate	5	<	Chlorobenzene	1	<	Nitrobenzene	5	<
cis-1,2,-Dichloroethene	1	<	1,1,1,2-Tetrachloroethane	1	<	1,2,4-Trichlorobenzene	1	<
Tetrahydrofuran	5	<	Ethylbenzene	1	<	Naphthalene	25	<
Methyacrylonitrile	5	<	m/p-xylene	1	<	1,2,3-Trichlorobenzene	1	<
Isobutyl alcohol	5	<	Bromoform	2	<	Hexachlorobutadiene	1	<
1,1,1-Trichloroethane	1	<	o-xylene	1	<			

Legend:
< = less than reporting limits
J = less than reporting limit, above detection levels
E = estimated, above calibration range

Surrogate Recovery	%
Dibromofluoromethane	84
Toluene-d8	98
4-Bromofluorobenzene	92

Notes:

Chemist/Reviewer: RMB

Date: 5/31/06

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Certificate of Analysis

Sample Identification	MW-4
Depth (ft)	
Sample Matrix	water
Units:	ug/L
% Solid	NA
Dilution	1
Sample Preservation	4 deg C.
Sample Volume	40 mL.
Date Sampled	5/26/2006
Date Analyzed	5/26/2006

Client:	Woodard & Curran, Inc.
Site:	GP Mill
Loc.:	Old Town, Maine
Project No.:	
Chemist:	D. M. Twomey

Analytical Method: USEPA SW-846 8260B
Instrument: Agilent 5890 GCD MS
Detectors: Mass spectrometer
Column: DB-5, 30 meter, .25 um
Conc. System: Tekmar 3000 Purge & Trap System
Trap: Vocarb 3000

Method:
USEPA SW846 METHOD 8260B

Compound	RL ug/L	Results	Compound	RL ug/L	Results	Compound	RL ug/L	Results
Chloromethane	1	<	1,2-Dichloroethane	1	<	Styrene	1	<
Dichlorodifluoromethane	1	<	1,1-Dichloropropene	1	<	trans-1,4-Dichloro-2-butene	1	<
Chloroethane	1	<	Carbon Tetrachloride	1	<	cis-1,4,-Dichloro-2-butene	1	<
Vinyl chloride	1	<	Benzene	1	<	1,1,2,2-Tetrachloroethane	1	<
Bromomethane	2	<	2-Chloroethanol	5	<	1,2,3-Trichloropropane	1	<
Trichlorofluoromethane	1	<	1,2-Dichloropropane	1	<	Isopropylbenzene	1	<
Diethyl ether	5	<	Trichloroethene	1	<	Bromobenzene	1	<
Acetonitrile	5	<	Bromodichloromethane	1	<	2-Chlorotoluene	2	<
Idomethane	5	<	Dibromomethane	1	<	4-Chlorotoluene	2	<
1,1-Dichloroethene	1	<	Methyl methacrylate	5	<	n-propylbenzene	1	<
Carbon Disulfide	2	<	1,4-Dioxane	25	<	1,3,5-Trimethylbenzene	2	<
Allyl chloride	5	<	cis-1,3-Dichloropropene	1	<	Pentachloroethane	1	<
Methylene chloride	10	<	trans-1,3-Dichloropropene	1	<	1,2,4-Trimethylbenzene	2	<
Acrylonitrile	5	<	2-Nitropropane	5	<	tert-butylbenzene	1	<
1,1,2 - Trichlorotrifluoroethane	1	<	Toluene	1	<	sec-butylbenzene	1	<
trans-1,2-Dichloroethene	1	<	1,1,2-Trichloroethane	1	<	1,3-Dichlorobenzene	2	<
MTBE	1	<	Ethylmethacrylate	5	<	1,4-Dichlorobenzene	2	<
Chloroform	1	<	1,3-Dichloropropane	1	<	p-iso-Propyltoluene	1	<
1,1-Dichloroethane	1	<	Dibromochloromethane	1	<	1,2-Dichlorobenzene	2	<
Chloroprene	2	<	1,2-Dibromomethane	1	<	n-butylbenzene	1	<
2,2-Dichloropropane	1	<	Tetrachloroethene	1	<	1,2-Bromo-3-chloropropane	1	<
Methyl acrylate	5	<	Chlorobenzene	1	<	Nitrobenzene	5	<
cis-1,2,-Dichloroethene	1	<	1,1,1,2-Tetrachloroethane	1	<	1,2,4-Trichlorobenzene	1	<
Tetrahydrofuran	5	<	Ethylbenzene	1	<	Naphthalene	25	<
Methyacrylonitrile	5	<	m/p-xylene	1	<	1,2,3-Trichlorobenzene	1	<
Isobutyl alcohol	5	<	Bromoform	2	<	Hexachlorobutadiene	1	<
1,1,1-Trichloroethane	1	<	o-xylene	1	<			

Legend:
< = less than reporting limits
J = less than reporting limit, above detection levels
E = estimated, above calibration range

Surrogate Recovery	%
Dibromofluoromethane	96
Toluene-d8	112
4-Bromofluorobenzene	98

Notes:

Chemist/Reviewer: RMB

Date: 5/31/06

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Sample Identification	SB-40
Depth (ft)	1.5-5.5'
Sample Matrix	water
Units:	ug/L
% Solid	NA
Dilution	1
Sample Preservation	4 deg C.
Sample Volume	40 mL.
Date Sampled	5/26/2005
Date Analyzed	5/26/2005

Client:	**Woodard & Curran, Inc.**
Site:	**GP Mill**
Loc.:	**Old Town, Maine**
Project No.:	
Chemist:	D. M. Twomey

Analytical Method: USEPA SW-846 8260B
Instrument: Agilent 5890 GCD MS
Detectors: Mass spectrometer
Column: DB-5, 30 meter, .25 um
Conc. System: Tekmar 3000 Purge & Trap System
Trap: Vocarb 3000

Method:
USEPA SW846 METHOD 8260B

Compound	RL ug/L	Results	Compound	RL ug/L	Results	Compound	RL ug/L	Results
Chloromethane	1	<	1,2-Dichloroethane	1	<	Styrene	1	<
Dichlorodifluoromethane	1	<	1,1-Dichloropropene	1	<	trans-1,4-Dichloro-2-butene	1	<
Chloroethane	1	<	Carbon Tetrachloride	1	<	cis-1,4,-Dichloro-2-butene	1	<
Vinyl chloride	1	<	Benzene	1	<	1,1,2,2-Tetrachloroethane	1	<
Bromomethane	2	<	2-Chloroethanol	5	<	1,2,3-Trichloropropane	1	<
Trichlorofluoromethane	1	<	1,2-Dichloropropane	1	<	Isopropylbenzene	1	<
Diethyl ether	5	<	Trichloroethene	1	<	Bromobenzene	1	<
Acetonitrile	5	<	Bromodichloromethane	1	<	2-Chlorotoluene	2	<
Idomethane	5	<	Dibromomethane	1	<	4-Chlorotoluene	2	<
1,1-Dichloroethene	1	<	Methyl methacrylate	5	<	n-propylbenzene	1	<
Carbon Disulfide	2	<	1,4-Dioxane	25	<	1,3,5-Trimethylbenzene	2	<
Allyl chloride	5	<	cis-1,3-Dichloropropene	1	<	Pentachloroethane	1	<
Methylene chloride	10	<	trans-1,3-Dichloropropene	1	<	1,2,4-Trimethylbenzene	2	<
Acrylonitrile	5	<	2-Nitropropane	5	<	tert-butylbenzene	1	<
1,1,2 - Trichlorotrifluoroethane	1	<	Toluene	1	<	sec-butylbenzene	1	<
trans-1,2-Dichloroethene	1	<	1,1,2-Trichloroethane	1	<	1,3-Dichlorobenzene	2	<
MTBE	1	<	Ethylmethacrylate	5	<	1,4-Dichlorobenzene	2	<
Chloroform	1	<	1,3-Dichloropropane	1	<	p-iso-Propyltoluene	1	<
1,1-Dichloroethane	1	<	Dibromochloromethane	1	<	1,2-Dichlorobenzene	2	<
Chloroprene	2	<	1,2-Dibromomethane	1	<	n-butylbenzene	1	<
2,2-Dichloropropane	1	<	Tetrachloroethene	1	<	1,2-Bromo-3-chloropropane	1	<
Methyl acrylate	5	<	Chlorobenzene	1	<	Nitrobenzene	5	<
cis-1,2,-Dichloroethene	1	<	1,1,1,2-Tetrachloroethane	1	<	1,2,4-Trichlorobenzene	1	<
Tetrahydrofuran	5	<	Ethylbenzene	1	<	Naphthalene	25	<
Methyacrylonitrile	5	<	m/p-xylene	1	<	1,2,3-Trichlorobenzene	1	<
Isobutyl alcohol	5	<	Bromoform	2	<	Hexachlorobutadiene	1	<
1,1,1-Trichloroethane	1	<	o-xylene	1	<			

Legend:
< = less than reporting limits
J = less than reporting limit, above detection levels
E = estimated, above calibration range

Surrogate Recovery	%
Dibromofluoromethane	90
Toluene-d8	100
4-Bromofluorobenzene	95

Notes:

Chemist/Reviewer: RMB

Date: 5/31/06

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Sample Identification	MW_5
Depth (ft)	4-11.5'
Sample Matrix	water
Units:	ug/L
% Solid	NA
Dilution	1
Sample Preservation	4 deg C.
Sample Volume	40 mL.
Date Sampled	5/26/2006
Date Analyzed	5/26/2006

Client:	Woodard & Curran, Inc.
Site:	GP Mill
Loc.:	Old Town, Maine
Project No.:	
Chemist:	D. M. Twomey

Analytical Method: USEPA SW-846 8260B
Instrument: Agilent 5890 GCD MS
Detectors: Mass spectrometer
Column: DB-5, 30 meter, .25 um
Conc. System: Tekmar 3000 Purge & Trap System
Trap: Vocarb 3000

Method:
USEPA SW846 METHOD 8260B

Compound	RL ug/L	Results	Compound	RL ug/L	Results	Compound	RL ug/L	Results
Chloromethane	1	<	1,2-Dichloroethane	1	<	Styrene	1	<
Dichlorodifluoromethane	1	<	1,1-Dichloropropene	1	<	trans-1,4-Dichloro-2-butene	1	<
Chloroethane	1	<	Carbon Tetrachloride	1	<	cis-1,4,-Dichloro-2-butene	1	<
Vinyl chloride	1	<	Benzene	1	<	1,1,2,2-Tetrachloroethane	1	<
Bromomethane	2	<	2-Chloroethanol	5	<	1,2,3-Trichloropropane	1	<
Trichlorofluoromethane	1	<	1,2-Dichloropropane	1	<	Isopropylbenzene	1	<
Diethyl ether	5	<	Trichloroethene	1	<	Bromobenzene	1	<
Acetonitrile	5	<	Bromodichloromethane	1	<	2-Chlorotoluene	2	<
Idomethane	5	<	Dibromomethane	1	<	4-Chlorotoluene	2	<
1,1-Dichloroethene	1	<	Methyl methacrylate	5	<	n-propylbenzene	1	<
Carbon Disulfide	2	<	1,4-Dioxane	25	<	1,3,5-Trimethylbenzene	2	<
Allyl chloride	5	<	cis-1,3-Dichloropropene	1	<	Pentachloroethane	1	<
Methylene chloride	10	<	trans-1,3-Dichloropropene	1	<	1,2,4-Trimethylbenzene	2	<
Acrylonitrile	5	<	2-Nitropropane	5	<	tert-butylbenzene	1	<
1,1,2 - Trichlorotrifluoroethane	1	<	Toluene	1	<	sec-butylbenzene	1	<
trans-1,2-Dichloroethene	1	<	1,1,2-Trichloroethane	1	<	1,3-Dichlorobenzene	2	<
MTBE	1	<	Ethylmethacrylate	5	<	1,4-Dichlorobenzene	2	<
Chloroform	1	<	1,3-Dichloropropane	1	<	p-iso-Propyltoluene	1	<
1,1-Dichloroethane	1	<	Dibromochloromethane	1	<	1,2-Dichlorobenzene	2	<
Chloroprene	2	<	1,2-Dibromomethane	1	<	n-butylbenzene	1	<
2,2-Dichloropropane	1	<	Tetrachloroethene	1	<	1,2-Bromo-3-chloropropane	1	<
Methyl acrylate	5	<	Chlorobenzene	1	<	Nitrobenzene	5	<
cis-1,2,-Dichloroethene	1	<	1,1,1,2-Tetrachloroethane	1	<	1,2,4-Trichlorobenzene	1	<
Tetrahydrofuran	5	<	Ethylbenzene	1	<	Naphthalene	25	<
Methyacrylonitrile	5	<	m/p-xylene	1	<	1,2,3-Trichlorobenzene	1	<
Isobutyl alcohol	5	<	Bromoform	2	<	Hexachlorobutadiene	1	<
1,1,1-Trichloroethane	1	<	o-xylene	1	<			

Legend:
< = less than reporting limits
J = less than reporting limit, above detection levels
E = estimated, above calibration range

Surrogate Recovery	%
Dibromofluoromethane	87
Toluene-d8	85
4-Bromofluorobenzene	102

Notes:

Chemist/Reviewer: RMB

Date: 5/31/06

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Sample Identification	MW-3
Depth (ft)	6-10'
Sample Matrix	water
Units:	ug/L
% Solid	NA
Dilution	1
Sample Preservation	4 deg C.
Sample Volume	40 mL.
Date Sampled	5/26/2006
Date Analyzed	5/26/2006

Client:	**Woodard & Curran, Inc.**
Site:	**GP Mill**
Loc.:	**Old Town, Maine**
Project No.:	
Chemist:	D. M. Twomey

Analytical Method: USEPA SW-846 8260B
Instrument: Agilent 5890 GCD MS
Detectors: Mass spectrometer
Column: DB-5, 30 meter, .25 um
Conc. System: Tekmar 3000 Purge & Trap System
Trap: Vocarb 3000

Method:
USEPA SW846 METHOD 8260B

Compound	RL ug/L	Results	Compound	RL ug/L	Results	Compound	RL ug/L	Results
Chloromethane	1	<	1,2-Dichloroethane	1	<	Styrene	1	<
Dichlorodifluoromethane	1	<	1,1-Dichloropropene	1	<	trans-1,4-Dichloro-2-butene	1	<
Chloroethane	1	<	Carbon Tetrachloride	1	<	cis-1,4,-Dichloro-2-butene	1	<
Vinyl chloride	1	<	Benzene	1	<	1,1,2,2-Tetrachloroethane	1	<
Bromomethane	2	<	2-Chloroethanol	5	<	1,2,3-Trichloropropane	1	<
Trichlorofluoromethane	1	<	1,2-Dichloropropane	1	<	Isopropylbenzene	1	<
Diethyl ether	5	<	Trichloroethene	1	<	Bromobenzene	1	<
Acetonitrile	5	<	Bromodichloromethane	1	<	2-Chlorotoluene	2	<
Idomethane	5	<	Dibromomethane	1	<	4-Chlorotoluene	2	<
1,1-Dichloroethene	1	<	Methyl methacrylate	5	<	n-propylbenzene	1	<
Carbon Disulfide	2	<	1,4-Dioxane	25	<	1,3,5-Trimethylbenzene	2	<
Allyl chloride	5	<	cis-1,3-Dichloropropene	1	<	Pentachloroethane	1	<
Methylene chloride	10	<	trans-1,3-Dichloropropene	1	<	1,2,4-Trimethylbenzene	2	<
Acrylonitrile	5	<	2-Nitropropane	5	<	tert-butylbenzene	1	<
1,1,2 - Trichlorotrifluoroethane	1	<	Toluene	1	<	sec-butylbenzene	1	<
trans-1,2-Dichloroethene	1	<	1,1,2-Trichloroethane	1	<	1,3-Dichlorobenzene	2	<
MTBE	1	<	Ethylmethacrylate	5	<	1,4-Dichlorobenzene	2	<
Chloroform	1	<	1,3-Dichloropropane	1	<	p-iso-Propyltoluene	1	<
1,1-Dichloroethane	1	<	Dibromochloromethane	1	<	1,2-Dichlorobenzene	2	<
Chloroprene	2	<	1,2-Dibromomethane	1	<	n-butylbenzene	1	<
2,2-Dichloropropane	1	<	Tetrachloroethene	1	<	1,2-Bromo-3-chloropropane	1	<
Methyl acrylate	5	<	Chlorobenzene	1	<	Nitrobenzene	5	<
cis-1,2,-Dichloroethene	1	<	1,1,1,2-Tetrachloroethane	1	<	1,2,4-Trichlorobenzene	1	<
Tetrahydrofuran	5	<	Ethylbenzene	1	<	Naphthalene	25	<
Methyacrylonitrile	5	<	m/p-xylene	1	<	1,2,3-Trichlorobenzene	1	<
Isobutyl alcohol	5	<	Bromoform	2	<	Hexachlorobutadiene	1	<
1,1,1-Trichloroethane	1	<	o-xylene	1	<	*Chlorodifluorethane***	*1*	**3.0**

Legend:
< = less than reporting limits
J = less than reporting limit, above detection levels
E = estimated, above calibration range

Surrogate Recovery	%
Dibromofluoromethane	82
Toluene-d8	98
4-Bromofluorobenzene	95

Notes:
** **- non-target compound identified.**

Chemist/Reviewer: RMB

Date: 5/31/06

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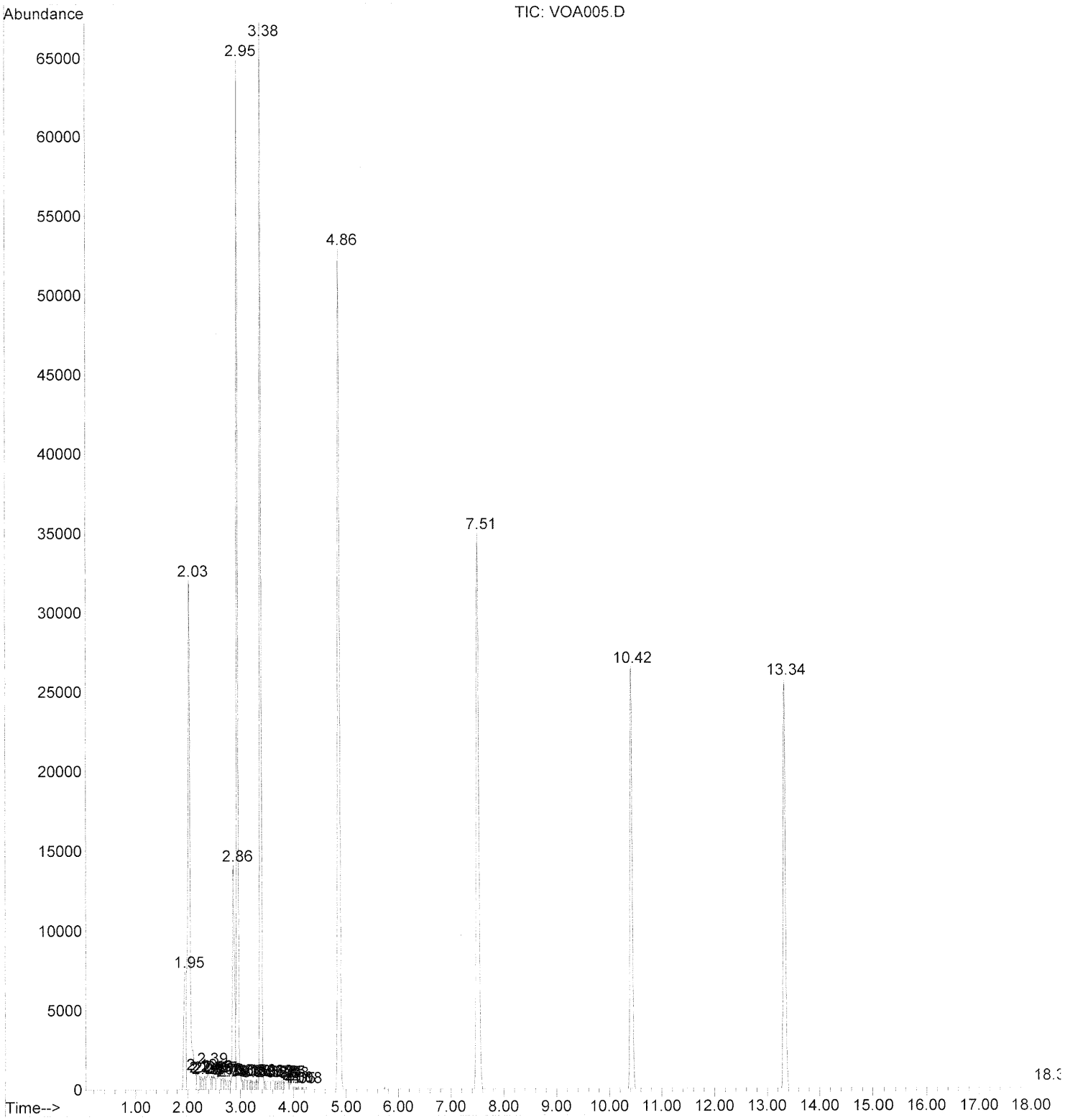
File : C:\HPCHEM\1\DATA\052406\VOA005.D
Operator : dmt
Acquired : 24 May 2006 12:21 pm using AcqMethod VOA0506
Instrument : GCD Syste
Sample Name: SB-11 6-9' H2O
Misc Info :
Vial Number: 1



Calibration Last Updated: Wed May 24 11:22:20 2006
Quantitation Report - Summary

Information from Data File:
File : C:\HPCHEM\1\DATA\052406\VOA005.D
Operator : dmt
Acquired : 24 May 2006 12:21 pm using AcqMethod VOA0506
Sample Name: SB-11 6-9' H2O
Misc Info :
Vial Number: 1
CurrentMeth: C:\HPCHEM\1\METHODS\VOA0506R.M

Quantitation Settings:
Non-Reference Peak Window: 10.00 Percent
Correlation Window: 0.02 Minutes
Default Multiplier: 1.00
Default Sample Concentration: 0.00
Peak Type Decoding: m -> Manually re-integrated

Number of Compounds in Database: 87

#	Compound	Ret Time	Concentration	Peak Type
1	Pentafluorobenzene	2.95	100.0 ppb	ISTD1
2	Dichlorofluoromethane	***** NOT FOUND *****		
3	Chloromethane	***** NOT FOUND *****		
4	Vinyl Chloride	***** NOT FOUND *****		
5	Bromomethane	***** NOT FOUND *****		
6	Chloroethane	***** NOT FOUND *****		
7	Trichlorofluoromethane	***** NOT FOUND *****		
8	Acetonitrile	***** NOT FOUND *****		
9	Diethyl ether	***** NOT FOUND *****		
10	1,1-Dichloroethene	***** NOT FOUND *****		
11	Acrylonitrile	***** NOT FOUND *****		
12	1,1,2-Trichlorotrifluoroethane	***** NOT FOUND *****		
13	Methylene Chloride	2.38	4.3 ppb	#
14	Iodomethane	***** NOT FOUND *****		
15	trans-1,2-Dichloroethene	***** NOT FOUND *****		
16	MTBE	***** NOT FOUND *****		
17	Allyl chloride	***** NOT FOUND *****		
18	1,1-Dichloroethane	***** NOT FOUND *****		
19	Carbon disulfide	***** NOT FOUND *****		
20	Chloroprene	***** NOT FOUND *****		
21	Methacrylonitrile	***** NOT FOUND *****		
22	cis-1,2-Dichloroethene	***** NOT FOUND *****		
23	Methyl acrylate	***** NOT FOUND *****		
24	Chloroform	***** NOT FOUND *****		
25	2,2-Dichloropropane	***** NOT FOUND *****		
26	Dibromofluoromethane-surrogate	2.86	92.3 ppb	
27	Isobutyl alcohol	***** NOT FOUND *****		
28	Tetrahydrofuran	***** NOT FOUND *****		
29	1,1,1-Trichloroethane	***** NOT FOUND *****		
30	1,2-Dichloroethane	***** NOT FOUND *****		
31	1,1-Dichloropropene	***** NOT FOUND *****		
32	Benzene	***** NOT FOUND *****		

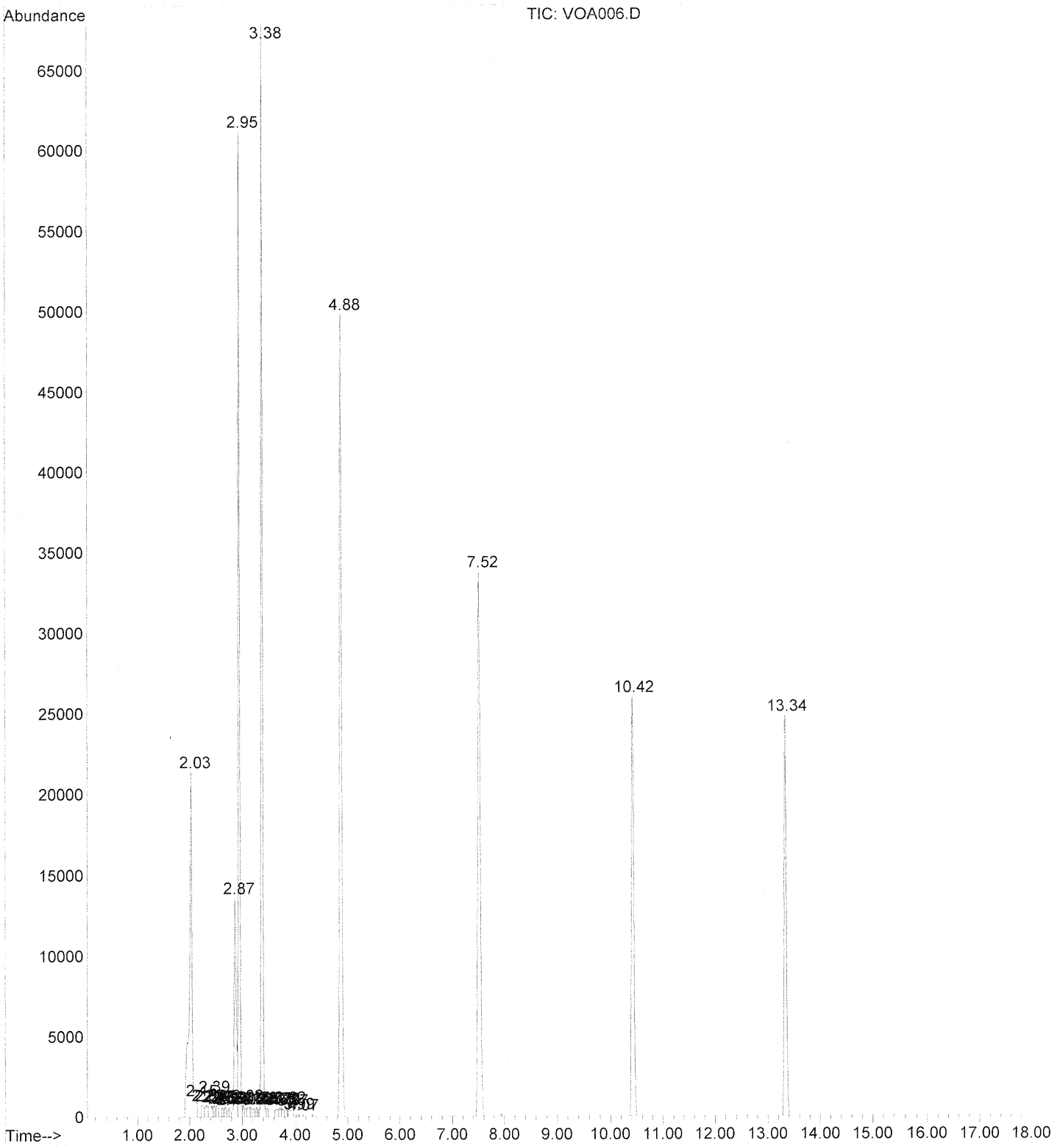
33 Carbon Tetrachloride	***** NOT FOUND *****			
34 2-Chloroethanol	***** NOT FOUND *****			
35 1,4-Difluorobenzene	3.38	100.0 ppb	ISTD2	
36 1,2-Dichloropropane	***** NOT FOUND *****			
37 Trichloroethene	***** NOT FOUND *****			
38 Dibromomethane	***** NOT FOUND *****			
39 Bromodichloromethane	***** NOT FOUND *****			
40 Methyl methacrylate	***** NOT FOUND *****			
41 1,4-Dioxane	***** NOT FOUND *****			
42 cis-1,3-Dichloropropene	***** NOT FOUND *****			
43 2-Nitropropane	***** NOT FOUND *****			
44 trans-1,3-Dichloropropene	***** NOT FOUND *****			
45 Toluene	***** NOT FOUND *****			
46 Toluene-d8-surrogate	4.86	95.9 ppb		96 % R
47 1,1,2-Trichloroethane	***** NOT FOUND *****			
48 1,3-Dichloropropane	***** NOT FOUND *****			
49 Ethylmethacrylate	***** NOT FOUND *****			
50 Dibromochloromethane	***** NOT FOUND *****			
51 1,2-Dibromomethane	***** NOT FOUND *****			
52 Tetrachloroethene	***** NOT FOUND *****			
53 Chlorobenzene	***** NOT FOUND *****			
54 Chlorobenzene-d5	7.50	100.0 ppb	ISTD3	
55 1,1,1,2-Tetrachloroethane	***** NOT FOUND *****			
56 Ethylbenzene	***** NOT FOUND *****			
57 m/p-xylene	***** NOT FOUND *****			
58 Bromoform	***** NOT FOUND *****			
59 Styrene	***** NOT FOUND *****			
60 o-Xylene	***** NOT FOUND *****			
61 cis-1,4-Dichloro-2-butene	***** NOT FOUND *****			
62 trans-1,4-Dichloro-2-butene	***** NOT FOUND *****			
63 Bromofluorobenzene-surrogate	10.42	96.8 ppb		97 % R
64 1,1,2,2-Tetrachloroethane	***** NOT FOUND *****			
65 1,2,3-Trichloropropane	***** NOT FOUND *****			
66 Isopropylbenzene	***** NOT FOUND *****			
67 Bromobenzene	***** NOT FOUND *****			
68 2-Chlorotoluene	***** NOT FOUND *****			
69 n-Propylbenzene	***** NOT FOUND *****			
70 4-Chlorotoluene	***** NOT FOUND *****			
71 1,3,5-Trimethylbenzene	***** NOT FOUND *****			
72 Pentachloroethane	***** NOT FOUND *****			
73 tert-Butylbenzene	***** NOT FOUND *****			
74 1,2,4-Trimethylbenzene	***** NOT FOUND *****			
75 1,3-Dichlorobenzene	***** NOT FOUND *****			
76 1,4-Dichlorobenzene-d4	13.34	100.0 ppb	ISTD4	
77 sec-Butylbenzene	***** NOT FOUND *****			
78 1,4-Dichlorobenzene	***** NOT FOUND *****			
79 p-iso-Propyltoluene	***** NOT FOUND *****			
80 1,2-Dichlorobenzene	***** NOT FOUND *****			
81 n-Butylbenzene	***** NOT FOUND *****			
82 1,2-Bromo-3-chloropropane	***** NOT FOUND *****			
83 Nitobenzene	***** NOT FOUND *****			
84 1,2,4-Trichlorobenzene	***** NOT FOUND *****			
85 Naphthalene	18.38	7.9 ppb	#	<
86 1,2,3-Trichlorobenzene	***** NOT FOUND *****			
87 Hexachlorobutadiene	***** NOT FOUND *****			

END OF REPORT

Qualifiers Not Satisfied

Wed May 24 12:40:26 2006

File : C:\HPCHEM\1\DATA\052406\VOA006.D
Operator : dmt
Acquired : 24 May 2006 12:47 pm using AcqMethod VOA0506
Instrument : GCD Syste
Sample Name: SB-12 6-9' H2O
Misc Info :
Vial Number: 1



Calibration Last Updated: Wed May 24 11:22:20 2006
Quantitation Report - Summary

Information from Data File:
File : C:\HPCHEM\1\DATA\052406\VOA006.D
Operator : dmt
Acquired : 24 May 2006 12:47 pm using AcqMethod VOA0506
Sample Name: SB-12 6-9' H2O
Misc Info :
Vial Number: 1
CurrentMeth: C:\HPCHEM\1\METHODS\VOA0506R.M

Quantitation Settings:
Non-Reference Peak Window: 10.00 Percent
Correlation Window: 0.02 Minutes
Default Multiplier: 1.00
Default Sample Concentration: 0.00
Peak Type Decoding: m -> Manually re-integrated

Number of Compounds in Database: 87

#	Compound	Ret Time	Concentration	Peak Type
1	Pentafluorobenzene	2.95	100.0 ppb	ISTD1
2	Dichlorofluoromethane	***** NOT FOUND *****		
3	Chloromethane	***** NOT FOUND *****		
4	Vinyl Chloride	***** NOT FOUND *****		
5	Bromomethane	***** NOT FOUND *****		
6	Chloroethane	***** NOT FOUND *****		
7	Trichlorofluoromethane	***** NOT FOUND *****		
8	Acetonitrile	***** NOT FOUND *****		
9	Diethyl ether	***** NOT FOUND *****		
10	1,1-Dichloroethene	***** NOT FOUND *****		
11	Acrylonitrile	***** NOT FOUND *****		
12	1,1,2-Trichlorotrifluoroethane	***** NOT FOUND *****		
13	Methylene Chloride	2.39	4.6 ppb	#
14	Iodomethane	***** NOT FOUND *****		
15	trans-1,2-Dichloroethene	***** NOT FOUND *****		
16	MTBE	***** NOT FOUND *****		
17	Allyl chloride	***** NOT FOUND *****		
18	1,1-Dichloroethane	***** NOT FOUND *****		
19	Carbon disulfide	***** NOT FOUND *****		
20	Chloroprene	***** NOT FOUND *****		
21	Methacrylonitrile	***** NOT FOUND *****		
22	cis-1,2-Dichloroethene	***** NOT FOUND *****		
23	Methyl acrylate	***** NOT FOUND *****		
24	Chloroform	***** NOT FOUND *****		
25	2,2-Dichloropropane	***** NOT FOUND *****		
26	Dibromofluoromethane-surrogate	2.87	92.2 ppb	
27	Isobutyl alcohol	***** NOT FOUND *****		
28	Tetrahydrofuran	***** NOT FOUND *****		
29	1,1,1-Trichloroethane	***** NOT FOUND *****		
30	1,2-Dichloroethane	***** NOT FOUND *****		
31	1,1-Dichloropropene	***** NOT FOUND *****		
32	Benzene	***** NOT FOUND *****		

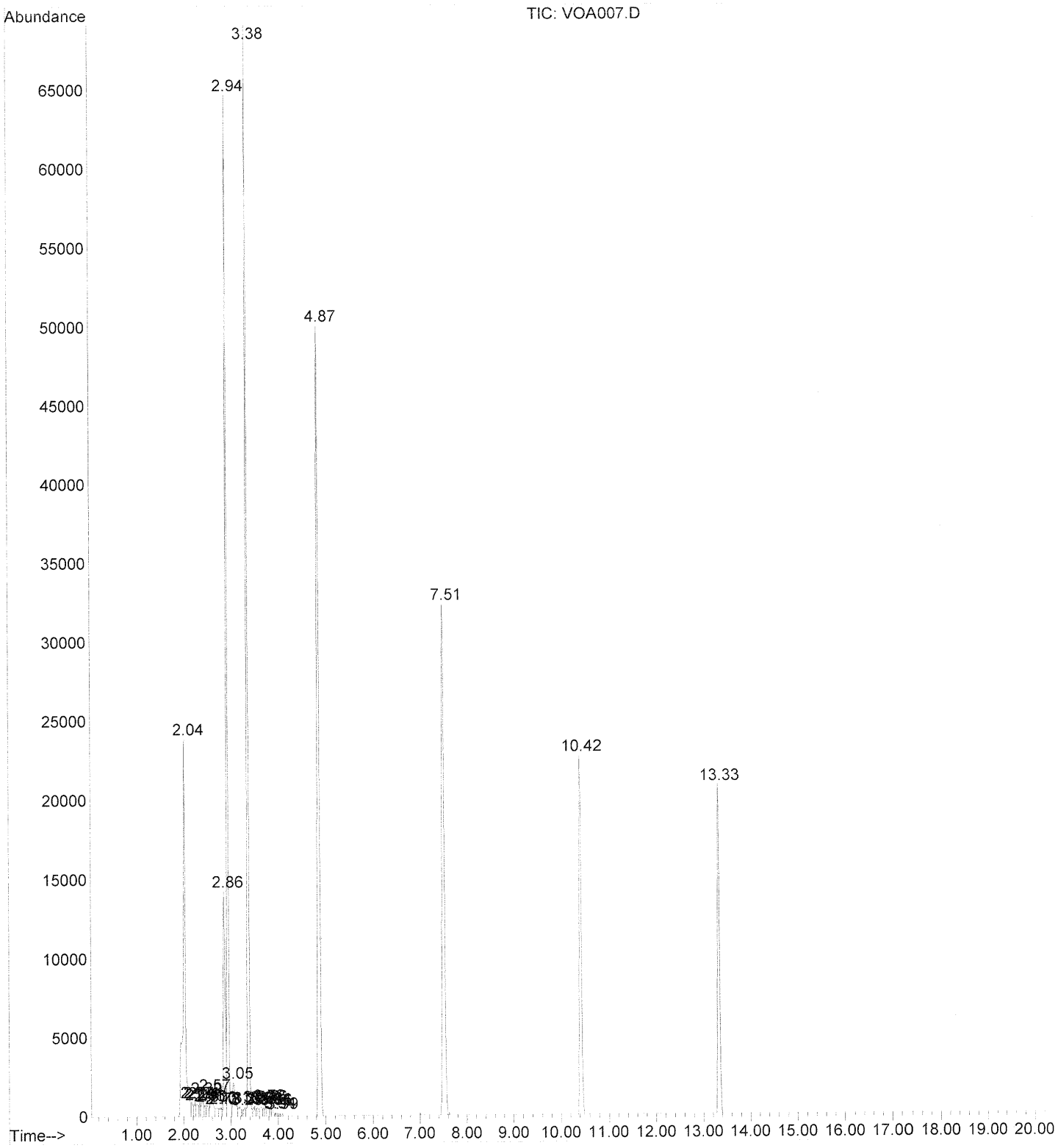
33 Carbon Tetrachloride	***** NOT FOUND *****			
34 2-Chloroethanol	***** NOT FOUND *****			
35 1,4-Difluorobenzene	3.38	100.0 ppb	ISTD2	
36 1,2-Dichloropropane	***** NOT FOUND *****			
37 Trichloroethene	***** NOT FOUND *****			
38 Dibromomethane	***** NOT FOUND *****			
39 Bromodichloromethane	***** NOT FOUND *****			
40 Methyl methacrylate	***** NOT FOUND *****			
41 1,4-Dioxane	***** NOT FOUND *****			
42 cis-1,3-Dichloropropene	***** NOT FOUND *****			
43 2-Nitropropane	***** NOT FOUND *****			
44 trans-1,3-Dichloropropene	***** NOT FOUND *****			
45 Toluene	***** NOT FOUND *****			95 %R
46 Toluene-d8-surrogate	4.87	94.7 ppb		
47 1,1,2-Trichloroethane	***** NOT FOUND *****			
48 1,3-Dichloropropane	***** NOT FOUND *****			
49 Ethylmethacrylate	***** NOT FOUND *****			
50 Dibromochloromethane	***** NOT FOUND *****			
51 1,2-Dibromomethane	***** NOT FOUND *****			
52 Tetrachloroethene	***** NOT FOUND *****			
53 Chlorobenzene	***** NOT FOUND *****			
54 Chlorobenzene-d5	7.52	100.0 ppb	ISTD3	
55 1,1,1,2-Tetrachloroethane	***** NOT FOUND *****			
56 Ethylbenzene	***** NOT FOUND *****			
57 m/p-xylene	***** NOT FOUND *****			
58 Bromoform	***** NOT FOUND *****			
59 Styrene	***** NOT FOUND *****			
60 o-Xylene	***** NOT FOUND *****			
61 cis-1,4-Dichloro-2-butene	***** NOT FOUND *****			
62 trans-1,4-Dichloro-2-butene	***** NOT FOUND *****			
63 Bromofluorobenzene-surrogate	10.42	90.0 ppb		90 %R
64 1,1,2,2-Tetrachloroethane	***** NOT FOUND *****			
65 1,2,3-Trichloropropane	***** NOT FOUND *****			
66 Isopropylbenzene	***** NOT FOUND *****			
67 Bromobenzene	***** NOT FOUND *****			
68 2-Chlorotoluene	***** NOT FOUND *****			
69 n-Propylbenzene	***** NOT FOUND *****			
70 4-Chlorotoluene	***** NOT FOUND *****			
71 1,3,5-Trimethylbenzene	***** NOT FOUND *****			
72 Pentachloroethane	***** NOT FOUND *****			
73 tert-Butylbenzene	***** NOT FOUND *****			
74 1,2,4-Trimethylbenzene	***** NOT FOUND *****			
75 1,3-Dichlorobenzene	***** NOT FOUND *****			
76 1,4-Dichlorobenzene-d4	13.34	100.0 ppb	ISTD4	
77 sec-Butylbenzene	***** NOT FOUND *****			
78 1,4-Dichlorobenzene	***** NOT FOUND *****			
79 p-iso-Propyltoluene	***** NOT FOUND *****			
80 1,2-Dichlorobenzene	***** NOT FOUND *****			
81 n-Butylbenzene	***** NOT FOUND *****			
82 1,2-Bromo-3-chloropropane	***** NOT FOUND *****			
83 Nitobenzene	***** NOT FOUND *****			
84 1,2,4-Trichlorobenzene	***** NOT FOUND *****			
85 Naphthalene	***** NOT FOUND *****			
86 1,2,3-Trichlorobenzene	***** NOT FOUND *****			
87 Hexachlorobutadiene	***** NOT FOUND *****			

END OF REPORT

Qualifiers Not Satisfied

Wed May 24 13:07:02 2006

File : C:\HPCHEM\1\DATA\052406\VOA007.D
Operator : dmt
Acquired : 24 May 2006 1:27 pm using AcqMethod VOA0506
Instrument : GCD Syste
Sample Name: SB-13 6-10' H2O
Misc Info :
Vial Number: 1



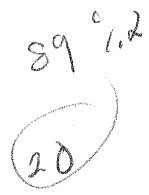
Calibration Last Updated: Wed May 24 11:22:20 2006
Quantitation Report - Summary

Information from Data File:
File : C:\HPCHEM\1\DATA\052406\VOA007.D
Operator : dmt
Acquired : 24 May 2006 1:27 pm using AcqMethod VOA0506
Sample Name: SB-13 6-10' H2O
Misc Info :
Vial Number: 1
CurrentMeth: C:\HPCHEM\1\METHODS\VOA0506R.M

Quantitation Settings:
Non-Reference Peak Window: 10.00 Percent
Correlation Window: 0.02 Minutes
Default Multiplier: 1.00
Default Sample Concentration: 0.00
Peak Type Decoding: m -> Manually re-integrated

Number of Compounds in Database: 87

#	Compound	Ret Time	Concentration	Peak Type
1	Pentafluorobenzene	2.94	100.0 ppb	ISTD1
2	Dichlorofluoromethane	***** NOT FOUND *****		
3	Chloromethane	***** NOT FOUND *****		
4	Vinyl Chloride	***** NOT FOUND *****		
5	Bromomethane	***** NOT FOUND *****		
6	Chloroethane	***** NOT FOUND *****		
7	Trichlorofluoromethane	***** NOT FOUND *****		
8	Acetonitrile	***** NOT FOUND *****		
9	Diethyl ether	***** NOT FOUND *****		
10	1,1-Dichloroethene	2.34	2.2 ppb	#
11	Acrylonitrile	***** NOT FOUND *****		
12	1,1,2-Trichlorotrifluoroethane	***** NOT FOUND *****		
13	Methylene Chloride	2.39	4.7 ppb	#
14	Iodomethane	***** NOT FOUND *****		
15	trans-1,2-Dichloroethene	***** NOT FOUND *****		
16	MTBE	***** NOT FOUND *****		
17	Allyl chloride	***** NOT FOUND *****		
18	1,1-Dichloroethane	2.58	40.1 ppb	#
19	Carbon disulfide	***** NOT FOUND *****		
20	Chloroprene	***** NOT FOUND *****		
21	Methacrylonitrile	***** NOT FOUND *****		
22	cis-1,2-Dichloroethene	***** NOT FOUND *****		
23	Methyl acrylate	***** NOT FOUND *****		
24	Chloroform	***** NOT FOUND *****		
25	2,2-Dichloropropane	***** NOT FOUND *****		
26	Dibromofluoromethane-surrogate	2.86	88.9 ppb	
27	Isobutyl alcohol	***** NOT FOUND *****		
28	Tetrahydrofuran	***** NOT FOUND *****		
29	1,1,1-Trichloroethane	3.05	20.1 ppb	
30	1,2-Dichloroethane	***** NOT FOUND *****		
31	1,1-Dichloropropene	***** NOT FOUND *****		
32	Benzene	***** NOT FOUND *****		



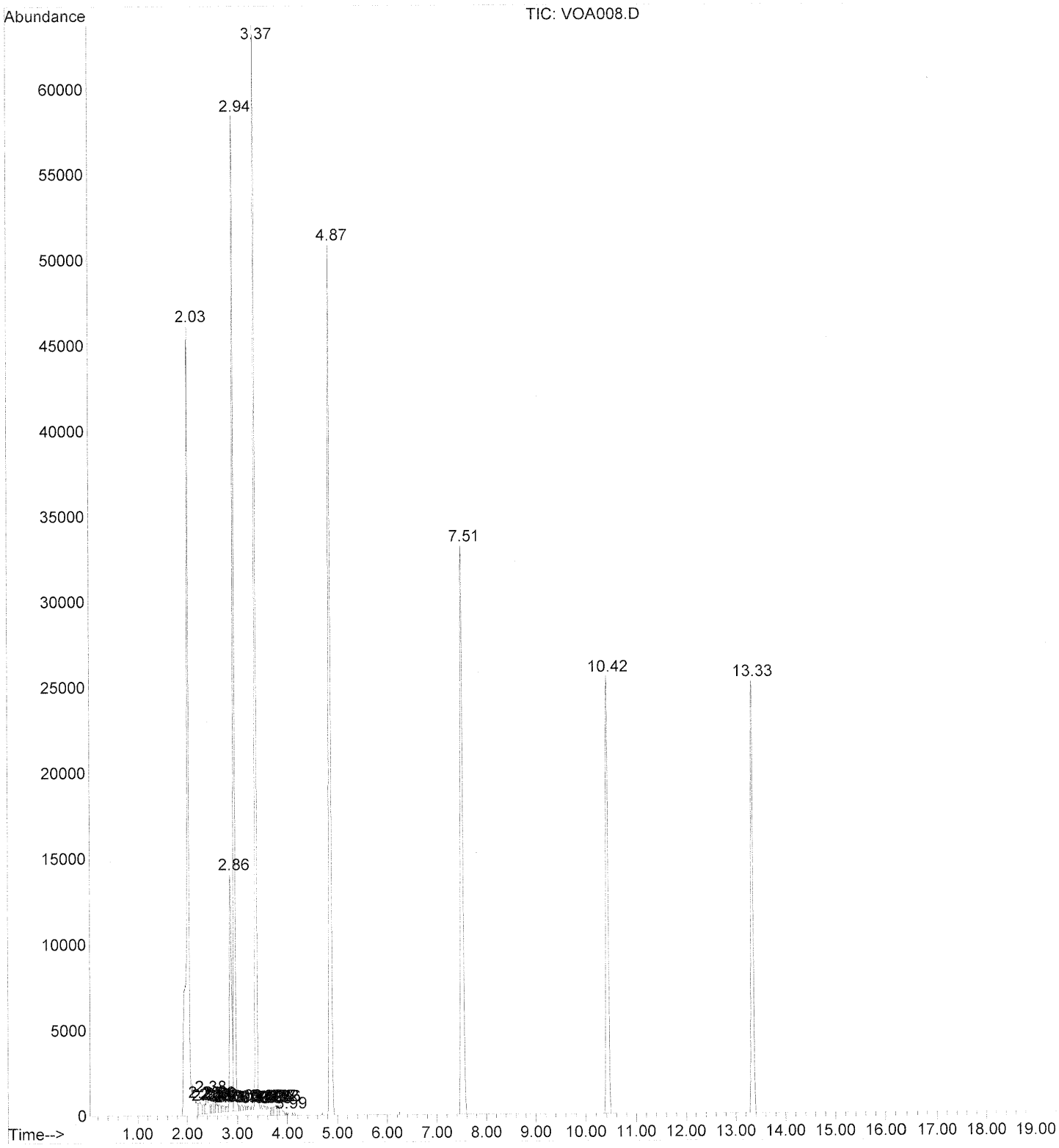
33 Carbon Tetrachloride	***** NOT FOUND *****			
34 2-Chloroethanol	***** NOT FOUND *****			
35 1,4-Difluorobenzene	3.38	100.0 ppb	ISTD2	
36 1,2-Dichloropropane	***** NOT FOUND *****			
37 Trichloroethene	***** NOT FOUND *****			
38 Dibromomethane	***** NOT FOUND *****			
39 Bromodichloromethane	***** NOT FOUND *****			
40 Methyl methacrylate	***** NOT FOUND *****			
41 1,4-Dioxane	***** NOT FOUND *****			
42 cis-1,3-Dichloropropene	***** NOT FOUND *****			
43 2-Nitropropane	***** NOT FOUND *****			
44 trans-1,3-Dichloropropene	***** NOT FOUND *****			
45 Toluene	***** NOT FOUND *****			
46 Toluene-d8-surrogate	4.87	90.9 ppb		91.8
47 1,1,2-Trichloroethane	***** NOT FOUND *****			
48 1,3-Dichloropropane	***** NOT FOUND *****			
49 Ethylmethacrylate	***** NOT FOUND *****			
50 Dibromochloromethane	***** NOT FOUND *****			
51 1,2-Dibromomethane	***** NOT FOUND *****			
52 Tetrachloroethene	***** NOT FOUND *****			
53 Chlorobenzene	***** NOT FOUND *****			
54 Chlorobenzene-d5	7.51	100.0 ppb	ISTD3	
55 1,1,1,2-Tetrachloroethane	***** NOT FOUND *****			
56 Ethylbenzene	***** NOT FOUND *****			
57 m/p-xylene	***** NOT FOUND *****			
58 Bromoform	***** NOT FOUND *****			
59 Styrene	***** NOT FOUND *****			
60 o-Xylene	***** NOT FOUND *****			
61 cis-1,4-Dichloro-2-butene	***** NOT FOUND *****			
62 trans-1,4-Dichloro-2-butene	***** NOT FOUND *****			
63 Bromofluorobenzene-surrogate	10.42	81.4 ppb		81.8
64 1,1,2,2-Tetrachloroethane	***** NOT FOUND *****			
65 1,2,3-Trichloropropane	***** NOT FOUND *****			
66 Isopropylbenzene	***** NOT FOUND *****			
67 Bromobenzene	***** NOT FOUND *****			
68 2-Chlorotoluene	***** NOT FOUND *****			
69 n-Propylbenzene	***** NOT FOUND *****			
70 4-Chlorotoluene	***** NOT FOUND *****			
71 1,3,5-Trimethylbenzene	***** NOT FOUND *****			
72 Pentachloroethane	***** NOT FOUND *****			
73 tert-Butylbenzene	***** NOT FOUND *****			
74 1,2,4-Trimethylbenzene	***** NOT FOUND *****			
75 1,3-Dichlorobenzene	***** NOT FOUND *****			
76 1,4-Dichlorobenzene-d4	13.33	100.0 ppb	ISTD4	
77 sec-Butylbenzene	***** NOT FOUND *****			
78 1,4-Dichlorobenzene	***** NOT FOUND *****			
79 p-iso-Propyltoluene	***** NOT FOUND *****			
80 1,2-Dichlorobenzene	***** NOT FOUND *****			
81 n-Butylbenzene	***** NOT FOUND *****			
82 1,2-Bromo-3-chloropropane	***** NOT FOUND *****			
83 Nitobenzene	***** NOT FOUND *****			
84 1,2,4-Trichlorobenzene	***** NOT FOUND *****			
85 Naphthalene	***** NOT FOUND *****			
86 1,2,3-Trichlorobenzene	***** NOT FOUND *****			
87 Hexachlorobutadiene	***** NOT FOUND *****			

END OF REPORT

Qualifiers Not Satisfied

Wed May 24 13:51:45 2006

File : C:\HPCHEM\1\DATA\052406\VOA008.D
Operator : dmt
Acquired : 24 May 2006 2:12 pm using AcqMethod VOA0506
Instrument : GCD Syste
Sample Name: SB-14 6-10' H2O
Misc Info :
Vial Number: 1



Calibration Last Updated: Wed May 24 11:22:20 2006
Quantitation Report - Summary

Information from Data File:
File : C:\HPCHEM\1\DATA\052406\VOA008.D
Operator : dmt
Acquired : 24 May 2006 2:12 pm using AcqMethod VOA0506
Sample Name: SB-14 6-10' H2O
Misc Info :
Vial Number: 1
CurrentMeth: C:\HPCHEM\1\METHODS\VOA0506R.M

Quantitation Settings:
Non-Reference Peak Window: 10.00 Percent
Correlation Window: 0.02 Minutes
Default Multiplier: 1.00
Default Sample Concentration: 0.00
Peak Type Decoding: m -> Manually re-integrated

Number of Compounds in Database: 87

#	Compound	Ret Time	Concentration	Peak Type
1	Pentafluorobenzene	2.95	100.0 ppb	ISTD1
2	Dichlorofluoromethane	***** NOT FOUND *****		
3	Chloromethane	***** NOT FOUND *****		
4	Vinyl Chloride	***** NOT FOUND *****		
5	Bromomethane	***** NOT FOUND *****		
6	Chloroethane	***** NOT FOUND *****		
7	Trichlorofluoromethane	***** NOT FOUND *****		
8	Acetonitrile	***** NOT FOUND *****		
9	Diethyl ether	***** NOT FOUND *****		
10	1,1-Dichloroethene	***** NOT FOUND *****		
11	Acrylonitrile	***** NOT FOUND *****		
12	1,1,2-Trichlorotrifluoroethane	***** NOT FOUND *****		
13	Methylene Chloride	2.38	3.7 ppb	#
14	Iodomethane	***** NOT FOUND *****		
15	trans-1,2-Dichloroethene	***** NOT FOUND *****		
16	MTBE	***** NOT FOUND *****		
17	Allyl chloride	***** NOT FOUND *****		
18	1,1-Dichloroethane	***** NOT FOUND *****		
19	Carbon disulfide	***** NOT FOUND *****		
20	Chloroprene	***** NOT FOUND *****		
21	Methacrylonitrile	***** NOT FOUND *****		
22	cis-1,2-Dichloroethene	***** NOT FOUND *****		
23	Methyl acrylate	***** NOT FOUND *****		
24	Chloroform	***** NOT FOUND *****		
25	2,2-Dichloropropane	***** NOT FOUND *****		
26	Dibromofluoromethane-surrogate	2.86	92.0 ppb	
27	Isobutyl alcohol	***** NOT FOUND *****		
28	Tetrahydrofuran	***** NOT FOUND *****		
29	1,1,1-Trichloroethane	***** NOT FOUND *****		
30	1,2-Dichloroethane	***** NOT FOUND *****		
31	1,1-Dichloropropene	***** NOT FOUND *****		
32	Benzene	***** NOT FOUND *****		

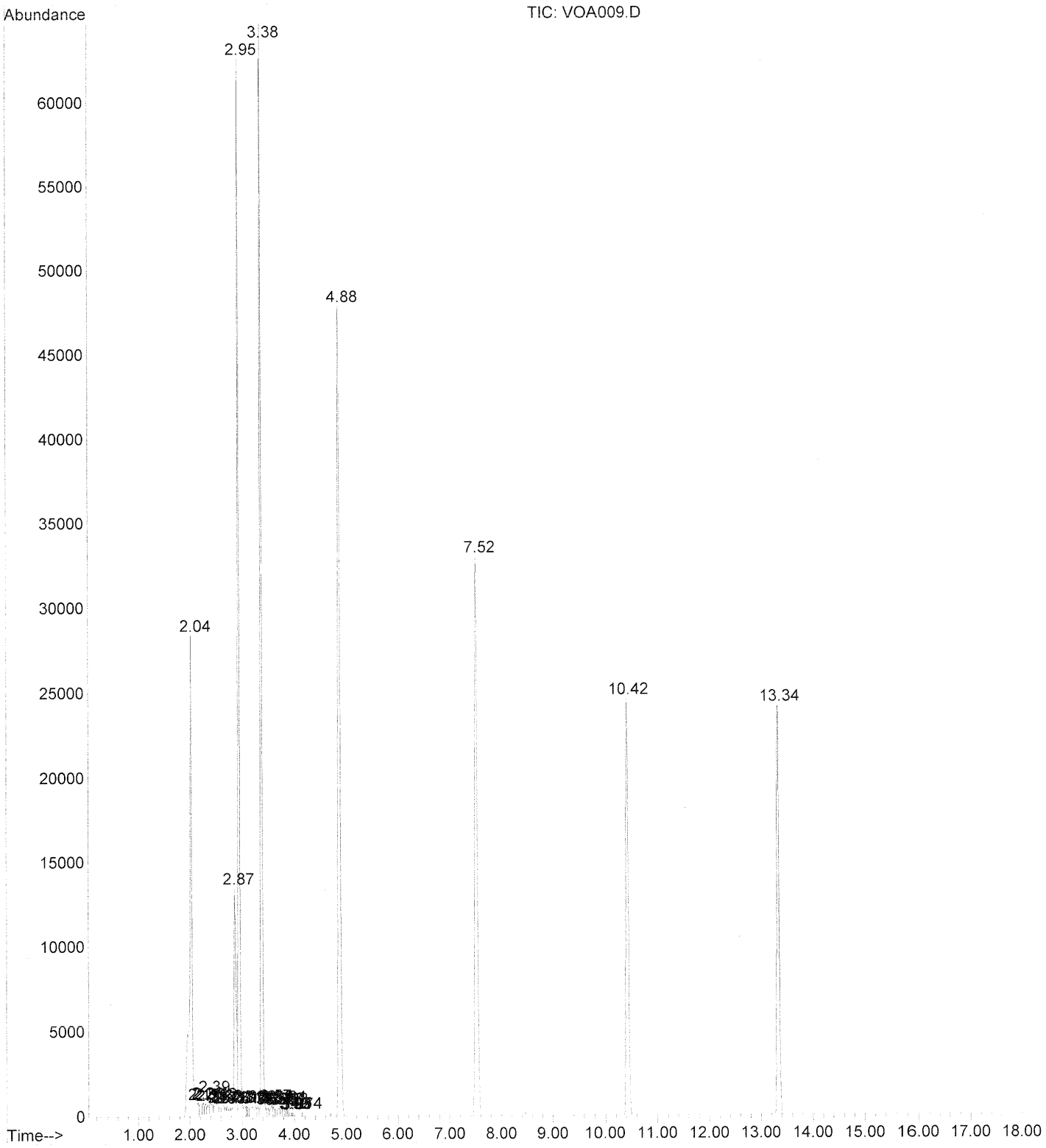
33	Carbon Tetrachloride	***** NOT FOUND *****		
34	2-Chloroethanol	***** NOT FOUND *****		
35	1,4-Difluorobenzene	3.37	100.0 ppb	ISTD2#
36	1,2-Dichloropropane	***** NOT FOUND *****		
37	Trichloroethene	***** NOT FOUND *****		
38	Dibromomethane	***** NOT FOUND *****		
39	Bromodichloromethane	***** NOT FOUND *****		
40	Methyl methacrylate	***** NOT FOUND *****		
41	1,4-Dioxane	***** NOT FOUND *****		
42	cis-1,3-Dichloropropene	***** NOT FOUND *****		
43	2-Nitropropane	***** NOT FOUND *****		
44	trans-1,3-Dichloropropene	***** NOT FOUND *****		
45	Toluene	***** NOT FOUND *****		
46	Toluene-d8-surrogate	4.87	94.9 ppb	
47	1,1,2-Trichloroethane	***** NOT FOUND *****		
48	1,3-Dichloropropane	***** NOT FOUND *****		
49	Ethylmethacrylate	***** NOT FOUND *****		
50	Dibromochloromethane	***** NOT FOUND *****		
51	1,2-Dibromomethane	***** NOT FOUND *****		
52	Tetrachloroethene	***** NOT FOUND *****		
53	Chlorobenzene	***** NOT FOUND *****		
54	Chlorobenzene-d5	7.52	100.0 ppb	ISTD3
55	1,1,1,2-Tetrachloroethane	***** NOT FOUND *****		
56	Ethylbenzene	***** NOT FOUND *****		
57	m/p-xylene	***** NOT FOUND *****		
58	Bromoform	***** NOT FOUND *****		
59	Styrene	***** NOT FOUND *****		
60	o-Xylene	***** NOT FOUND *****		
61	cis-1,4-Dichloro-2-butene	***** NOT FOUND *****		
62	trans-1,4-Dichloro-2-butene	***** NOT FOUND *****		
63	Bromofluorobenzene-surrogate	10.41	94.4 ppb	
64	1,1,2,2-Tetrachloroethane	***** NOT FOUND *****		
65	1,2,3-Trichloropropane	***** NOT FOUND *****		
66	Isopropylbenzene	***** NOT FOUND *****		
67	Bromobenzene	***** NOT FOUND *****		
68	2-Chlorotoluene	***** NOT FOUND *****		
69	n-Propylbenzene	***** NOT FOUND *****		
70	4-Chlorotoluene	***** NOT FOUND *****		
71	1,3,5-Trimethylbenzene	***** NOT FOUND *****		
72	Pentachloroethane	***** NOT FOUND *****		
73	tert-Butylbenzene	***** NOT FOUND *****		
74	1,2,4-Trimethylbenzene	***** NOT FOUND *****		
75	1,3-Dichlorobenzene	***** NOT FOUND *****		
76	1,4-Dichlorobenzene-d4	13.34	100.0 ppb	ISTD4
77	sec-Butylbenzene	***** NOT FOUND *****		
78	1,4-Dichlorobenzene	***** NOT FOUND *****		
79	p-iso-Propyltoluene	***** NOT FOUND *****		
80	1,2-Dichlorobenzene	***** NOT FOUND *****		
81	n-Butylbenzene	***** NOT FOUND *****		
82	1,2-Bromo-3-chloropropane	***** NOT FOUND *****		
83	Nitobenzene	***** NOT FOUND *****		
84	1,2,4-Trichlorobenzene	***** NOT FOUND *****		
85	Naphthalene	***** NOT FOUND *****		
86	1,2,3-Trichlorobenzene	***** NOT FOUND *****		
87	Hexachlorobutadiene	***** NOT FOUND *****		

END OF REPORT

Qualifiers Not Satisfied

Wed May 24 14:33:05 2006

File : C:\HPCHEM\1\DATA\052406\VOA009.D
Operator : dmt
Acquired : 24 May 2006 2:44 pm using AcqMethod VOA0506
Instrument : GCD Syste
Sample Name: ~~SB-16~~ SB-17 3.5-7.5' H2O
Misc Info : DUP
Vial Number: 1



Calibration Last Updated: Wed May 24 11:22:20 2006
Quantitation Report - Summary

Information from Data File:
File : C:\HPCHEM\1\DATA\052406\VOA009.D
Operator : dmt
Acquired : 24 May 2006 2:44 pm using AcqMethod VOA0506
Sample Name: ~~SB-16~~ 3.5-7.5' H2O
Misc Info : SB-17
Vial Number: 1 (DUP)
CurrentMeth: C:\HPCHEM\1\METHODS\VOA0506R.M

Quantitation Settings:
Non-Reference Peak Window: 10.00 Percent
Correlation Window: 0.02 Minutes
Default Multiplier: 1.00
Default Sample Concentration: 0.00
Peak Type Decoding: m -> Manually re-integrated

Number of Compounds in Database: 87

#	Compound	Ret Time	Concentration	Peak Type
1	Pentafluorobenzene	2.95	100.0 ppb	ISTD1
2	Dichlorofluoromethane	***** NOT FOUND *****		
3	Chloromethane	***** NOT FOUND *****		
4	Vinyl Chloride	***** NOT FOUND *****		
5	Bromomethane	***** NOT FOUND *****		
6	Chloroethane	***** NOT FOUND *****		
7	Trichlorofluoromethane	***** NOT FOUND *****		
8	Acetonitrile	***** NOT FOUND *****		
9	Diethyl ether	***** NOT FOUND *****		
10	1,1-Dichloroethene	***** NOT FOUND *****		
11	Acrylonitrile	***** NOT FOUND *****		
12	1,1,2-Trichlorotrifluoroethane	***** NOT FOUND *****		
13	Methylene Chloride	2.39	3.3 ppb	#
14	Iodomethane	***** NOT FOUND *****		
15	trans-1,2-Dichloroethene	***** NOT FOUND *****		
16	MTBE	***** NOT FOUND *****		
17	Allyl chloride	***** NOT FOUND *****		
18	1,1-Dichloroethane	***** NOT FOUND *****		
19	Carbon disulfide	***** NOT FOUND *****		
20	Chloroprene	***** NOT FOUND *****		
21	Methacrylonitrile	***** NOT FOUND *****		
22	cis-1,2-Dichloroethene	***** NOT FOUND *****		
23	Methyl acrylate	***** NOT FOUND *****		
24	Chloroform	***** NOT FOUND *****		
25	2,2-Dichloropropane	***** NOT FOUND *****		
26	Dibromofluoromethane-surrogate	2.87	86.4 ppb	86
27	Isobutyl alcohol	***** NOT FOUND *****		
28	Tetrahydrofuran	***** NOT FOUND *****		
29	1,1,1-Trichloroethane	***** NOT FOUND *****		
30	1,2-Dichloroethane	***** NOT FOUND *****		
31	1,1-Dichloropropene	***** NOT FOUND *****		
32	Benzene	***** NOT FOUND *****		

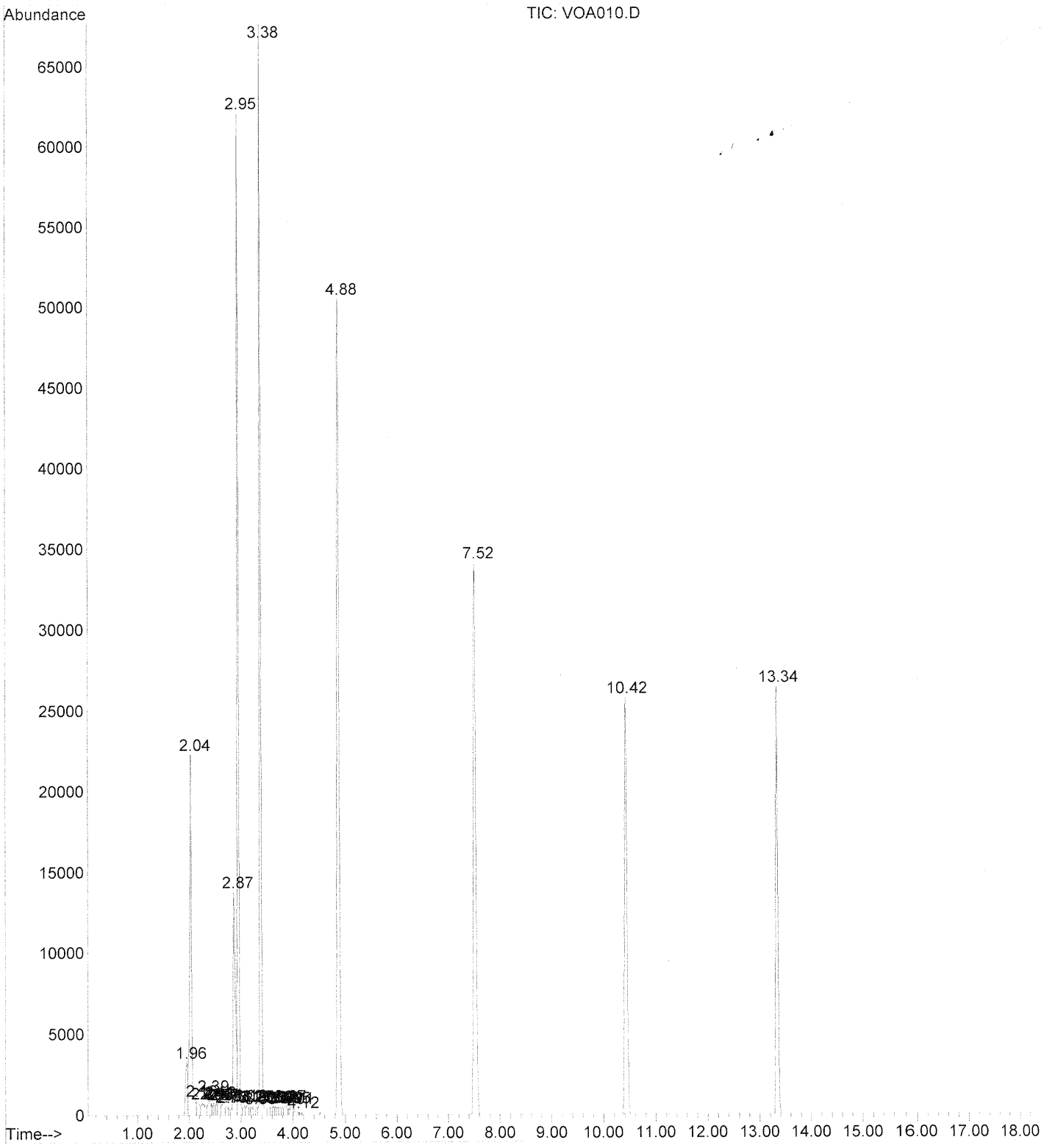
33 Carbon Tetrachloride	***** NOT FOUND *****			
34 2-Chloroethanol	***** NOT FOUND *****			
35 1,4-Difluorobenzene	3.38	100.0 ppb	ISTD2	
36 1,2-Dichloropropane	***** NOT FOUND *****			
37 Trichloroethene	***** NOT FOUND *****			
38 Dibromomethane	***** NOT FOUND *****			
39 Bromodichloromethane	***** NOT FOUND *****			
40 Methyl methacrylate	***** NOT FOUND *****			
41 1,4-Dioxane	***** NOT FOUND *****			
42 cis-1,3-Dichloropropene	***** NOT FOUND *****			
43 2-Nitropropane	***** NOT FOUND *****			
44 trans-1,3-Dichloropropene	***** NOT FOUND *****			
45 Toluene	***** NOT FOUND *****			
46 Toluene-d8-surrogate	4.88	95.7 ppb		96 %
47 1,1,2-Trichloroethane	***** NOT FOUND *****			
48 1,3-Dichloropropane	***** NOT FOUND *****			
49 Ethylmethacrylate	***** NOT FOUND *****			
50 Dibromochloromethane	***** NOT FOUND *****			
51 1,2-Dibromomethane	***** NOT FOUND *****			
52 Tetrachloroethene	***** NOT FOUND *****			
53 Chlorobenzene	***** NOT FOUND *****			
54 Chlorobenzene-d5	7.52	100.0 ppb	ISTD3	
55 1,1,1,2-Tetrachloroethane	***** NOT FOUND *****			
56 Ethylbenzene	***** NOT FOUND *****			
57 m/p-xylene	***** NOT FOUND *****			
58 Bromoform	***** NOT FOUND *****			
59 Styrene	***** NOT FOUND *****			
60 o-Xylene	***** NOT FOUND *****			
61 cis-1,4-Dichloro-2-butene	***** NOT FOUND *****			
62 trans-1,4-Dichloro-2-butene	***** NOT FOUND *****			
63 Bromofluorobenzene-surrogate	10.43	92.0 ppb		92 %
64 1,1,2,2-Tetrachloroethane	***** NOT FOUND *****			
65 1,2,3-Trichloropropane	***** NOT FOUND *****			
66 Isopropylbenzene	***** NOT FOUND *****			
67 Bromobenzene	***** NOT FOUND *****			
68 2-Chlorotoluene	***** NOT FOUND *****			
69 n-Propylbenzene	***** NOT FOUND *****			
70 4-Chlorotoluene	***** NOT FOUND *****			
71 1,3,5-Trimethylbenzene	***** NOT FOUND *****			
72 Pentachloroethane	***** NOT FOUND *****			
73 tert-Butylbenzene	***** NOT FOUND *****			
74 1,2,4-Trimethylbenzene	***** NOT FOUND *****			
75 1,3-Dichlorobenzene	***** NOT FOUND *****			
76 1,4-Dichlorobenzene-d4	13.33	100.0 ppb	ISTD4	
77 sec-Butylbenzene	***** NOT FOUND *****			
78 1,4-Dichlorobenzene	***** NOT FOUND *****			
79 p-iso-Propyltoluene	***** NOT FOUND *****			
80 1,2-Dichlorobenzene	***** NOT FOUND *****			
81 n-Butylbenzene	***** NOT FOUND *****			
82 1,2-Bromo-3-chloropropane	***** NOT FOUND *****			
83 Nitobenzene	***** NOT FOUND *****			
84 1,2,4-Trichlorobenzene	***** NOT FOUND *****			
85 Naphthalene	***** NOT FOUND *****			
86 1,2,3-Trichlorobenzene	***** NOT FOUND *****			
87 Hexachlorobutadiene	***** NOT FOUND *****			

END OF REPORT

Qualifiers Not Satisfied

Wed May 24 15:03:59 2006

File : C:\HPCHEM\1\DATA\052406\VOA010.D
Operator : dmt
Acquired : 24 May 2006 3:12 pm using AcqMethod VOA0506
Instrument : GCD Syste
Sample Name: ~~SB-18~~ 6-10' H2O
Misc Info : SB-15 dup
Vial Number: 1



Calibration Last Updated: Wed May 24 11:22:20 2006
Quantitation Report - Summary

Information from Data File:
File : C:\HPCHEM\1\DATA\052406\VOA010.D
Operator : dmt
Acquired : 24 May 2006 3:12 pm using AcqMethod VOA0506
Sample Name: ~~SB-18~~ 6-10' H2O
Misc Info : SB-15
Vial Number: 1
CurrentMeth: C:\HPCHEM\1\METHODS\VOA0506R.M

Quantitation Settings:
Non-Reference Peak Window: 10.00 Percent
Correlation Window: 0.02 Minutes
Default Multiplier: 1.00
Default Sample Concentration: 0.00
Peak Type Decoding: m -> Manually re-integrated

Number of Compounds in Database: 87

#	Compound	Ret Time	Concentration	Peak Type
1	Pentafluorobenzene	2.95	100.0 ppb	ISTD1
2	Dichlorofluoromethane	***** NOT FOUND *****		
3	Chloromethane	***** NOT FOUND *****		
4	Vinyl Chloride	***** NOT FOUND *****		
5	Bromomethane	***** NOT FOUND *****		
6	Chloroethane	***** NOT FOUND *****		
7	Trichlorofluoromethane	***** NOT FOUND *****		
8	Acetonitrile	***** NOT FOUND *****		
9	Diethyl ether	***** NOT FOUND *****		
10	1,1-Dichloroethene	***** NOT FOUND *****		
11	Acrylonitrile	***** NOT FOUND *****		
12	1,1,2-Trichlorotrifluoroethane	***** NOT FOUND *****		
13	Methylene Chloride	2.39	4.2 ppb	#
14	Iodomethane	***** NOT FOUND *****		
15	trans-1,2-Dichloroethene	***** NOT FOUND *****		
16	MTBE	***** NOT FOUND *****		
17	Allyl chloride	***** NOT FOUND *****		
18	1,1-Dichloroethane	***** NOT FOUND *****		
19	Carbon disulfide	***** NOT FOUND *****		
20	Chloroprene	***** NOT FOUND *****		
21	Methacrylonitrile	***** NOT FOUND *****		
22	cis-1,2-Dichloroethene	***** NOT FOUND *****		
23	Methyl acrylate	***** NOT FOUND *****		
24	Chloroform	***** NOT FOUND *****		
25	2,2-Dichloropropane	***** NOT FOUND *****		
26	Dibromofluoromethane-surrogate	2.87	92.7 ppb	93 % R
27	Isobutyl alcohol	***** NOT FOUND *****		
28	Tetrahydrofuran	***** NOT FOUND *****		
29	1,1,1-Trichloroethane	***** NOT FOUND *****		
30	1,2-Dichloroethane	***** NOT FOUND *****		
31	1,1-Dichloropropene	***** NOT FOUND *****		
32	Benzene	***** NOT FOUND *****		

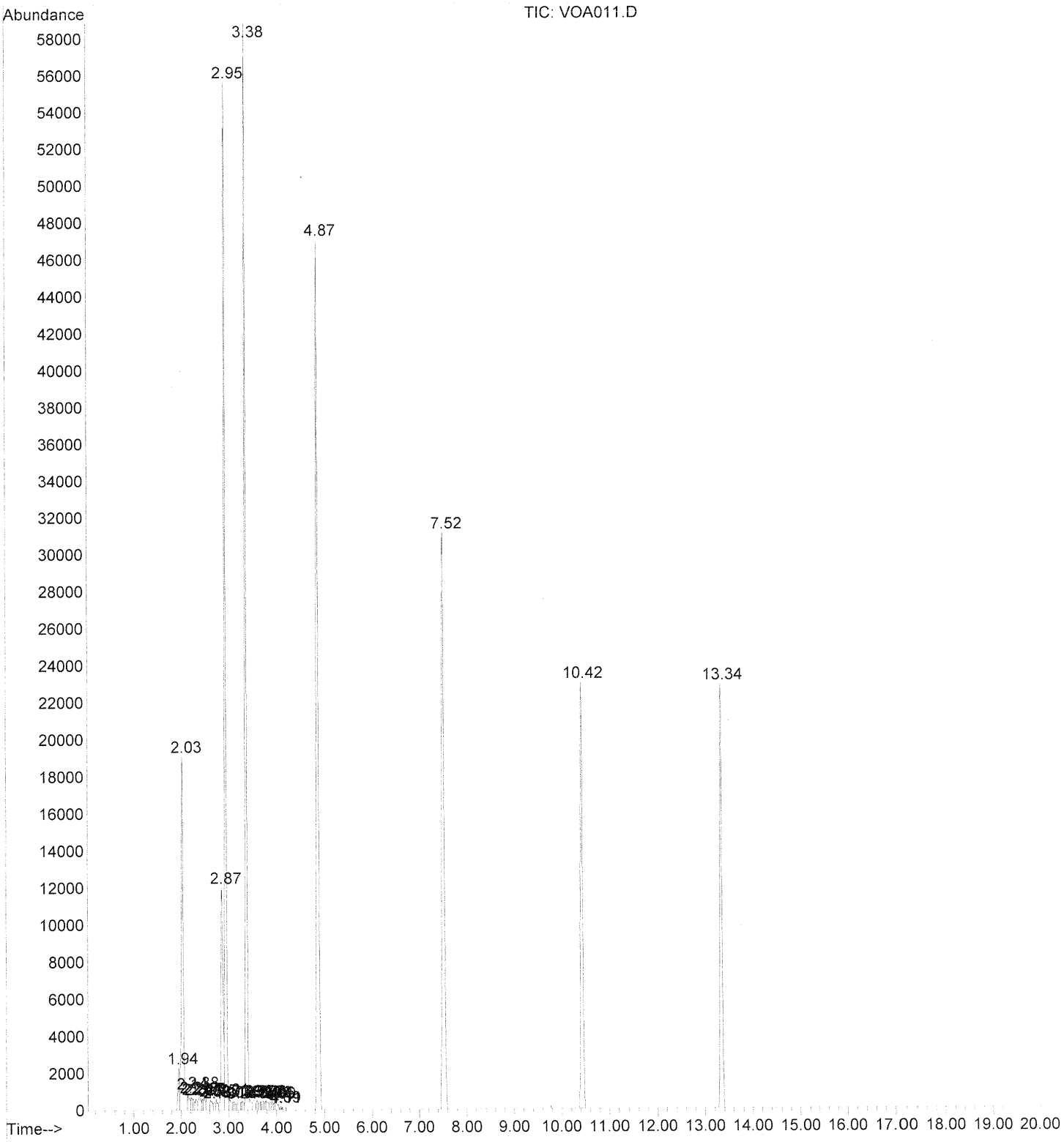
33 Carbon Tetrachloride	***** NOT FOUND *****			
34 2-Chloroethanol	***** NOT FOUND *****			
35 1,4-Difluorobenzene	3.38	100.0 ppb	ISTD2#	
36 1,2-Dichloropropane	***** NOT FOUND *****			
37 Trichloroethene	***** NOT FOUND *****			
38 Dibromomethane	***** NOT FOUND *****			
39 Bromodichloromethane	***** NOT FOUND *****			
40 Methyl methacrylate	***** NOT FOUND *****			
41 1,4-Dioxane	***** NOT FOUND *****			
42 cis-1,3-Dichloropropene	***** NOT FOUND *****			
43 2-Nitropropane	***** NOT FOUND *****			
44 trans-1,3-Dichloropropene	***** NOT FOUND *****			
45 Toluene	***** NOT FOUND *****			
46 Toluene-d8-surrogate	4.88	91.6 ppb	92 %R	
47 1,1,2-Trichloroethane	***** NOT FOUND *****			
48 1,3-Dichloropropane	***** NOT FOUND *****			
49 Ethylmethacrylate	***** NOT FOUND *****			
50 Dibromochloromethane	***** NOT FOUND *****			
51 1,2-Dibromomethane	***** NOT FOUND *****			
52 Tetrachloroethene	***** NOT FOUND *****			
53 Chlorobenzene	***** NOT FOUND *****			
54 Chlorobenzene-d5	7.52	100.0 ppb	ISTD3	
55 1,1,1,2-Tetrachloroethane	***** NOT FOUND *****			
56 Ethylbenzene	***** NOT FOUND *****			
57 m/p-xylene	***** NOT FOUND *****			
58 Bromoform	***** NOT FOUND *****			
59 Styrene	***** NOT FOUND *****			
60 o-Xylene	***** NOT FOUND *****			
61 cis-1,4-Dichloro-2-butene	***** NOT FOUND *****			
62 trans-1,4-Dichloro-2-butene	***** NOT FOUND *****			
63 Bromofluorobenzene-surrogate	10.42	91.1 ppb	91 %R	
64 1,1,2,2-Tetrachloroethane	***** NOT FOUND *****			
65 1,2,3-Trichloropropane	***** NOT FOUND *****			
66 Isopropylbenzene	***** NOT FOUND *****			
67 Bromobenzene	***** NOT FOUND *****			
68 2-Chlorotoluene	***** NOT FOUND *****			
69 n-Propylbenzene	***** NOT FOUND *****			
70 4-Chlorotoluene	***** NOT FOUND *****			
71 1,3,5-Trimethylbenzene	***** NOT FOUND *****			
72 Pentachloroethane	***** NOT FOUND *****			
73 tert-Butylbenzene	***** NOT FOUND *****			
74 1,2,4-Trimethylbenzene	***** NOT FOUND *****			
75 1,3-Dichlorobenzene	***** NOT FOUND *****			
76 1,4-Dichlorobenzene-d4	13.33	100.0 ppb	ISTD4	
77 sec-Butylbenzene	***** NOT FOUND *****			
78 1,4-Dichlorobenzene	***** NOT FOUND *****			
79 p-iso-Propyltoluene	***** NOT FOUND *****			
80 1,2-Dichlorobenzene	***** NOT FOUND *****			
81 n-Butylbenzene	***** NOT FOUND *****			
82 1,2-Bromo-3-chloropropane	***** NOT FOUND *****			
83 Nitobenzene	***** NOT FOUND *****			
84 1,2,4-Trichlorobenzene	***** NOT FOUND *****			
85 Naphthalene	***** NOT FOUND *****			
86 1,2,3-Trichlorobenzene	***** NOT FOUND *****			
87 Hexachlorobutadiene	***** NOT FOUND *****			

END OF REPORT

Qualifiers Not Satisfied

Wed May 24 15:32:22 2006

File : C:\HPCHEM\1\DATA\052406\VOA011.D
Operator : dmt
Acquired : 24 May 2006 3:39 pm using AcqMethod VOA0506
Instrument : GCD Syste
Sample Name: SB-19 2-6' H2O
Misc Info :
Vial Number: 1



Calibration Last Updated: Wed May 24 11:22:20 2006
Quantitation Report - Summary

Information from Data File:
File : C:\HPCHEM\1\DATA\052406\VOA011.D
Operator : dmt
Acquired : 24 May 2006 3:39 pm using AcqMethod VOA0506
Sample Name: SB-19 2-6' H2O
Misc Info :
Vial Number: 1
CurrentMeth: C:\HPCHEM\1\METHODS\VOA0506R.M

Quantitation Settings:
Non-Reference Peak Window: 10.00 Percent
Correlation Window: 0.02 Minutes
Default Multiplier: 1.00
Default Sample Concentration: 0.00
Peak Type Decoding: m -> Manually re-integrated

Number of Compounds in Database: 87

#	Compound	Ret Time	Concentration	Peak Type
1	Pentafluorobenzene	2.95	100.0 ppb	ISTD1
2	Dichlorofluoromethane	***** NOT FOUND *****		
3	Chloromethane	***** NOT FOUND *****		
4	Vinyl Chloride	***** NOT FOUND *****		
5	Bromomethane	***** NOT FOUND *****		
6	Chloroethane	***** NOT FOUND *****		
7	Trichlorofluoromethane	***** NOT FOUND *****		
8	Acetonitrile	***** NOT FOUND *****		
9	Diethyl ether	***** NOT FOUND *****		
10	1,1-Dichloroethene	***** NOT FOUND *****		
11	Acrylonitrile	***** NOT FOUND *****		
12	1,1,2-Trichlorotrifluoroethane	***** NOT FOUND *****		
13	Methylene Chloride	2.38	4.1 ppb	#
14	Iodomethane	***** NOT FOUND *****		
15	trans-1,2-Dichloroethene	***** NOT FOUND *****		
16	MTBE	***** NOT FOUND *****		
17	Allyl chloride	***** NOT FOUND *****		
18	1,1-Dichloroethane	***** NOT FOUND *****		
19	Carbon disulfide	***** NOT FOUND *****		
20	Chloroprene	***** NOT FOUND *****		
21	Methacrylonitrile	***** NOT FOUND *****		
22	cis-1,2-Dichloroethene	***** NOT FOUND *****		
23	Methyl acrylate	***** NOT FOUND *****		
24	Chloroform	***** NOT FOUND *****		
25	2,2-Dichloropropane	***** NOT FOUND *****		
26	Dibromofluoromethane-surrogate	2.87	91.3 ppb	
27	Isobutyl alcohol	***** NOT FOUND *****		
28	Tetrahydrofuran	***** NOT FOUND *****		
29	1,1,1-Trichloroethane	***** NOT FOUND *****		
30	1,2-Dichloroethane	***** NOT FOUND *****		
31	1,1-Dichloropropene	***** NOT FOUND *****		
32	Benzene	***** NOT FOUND *****		

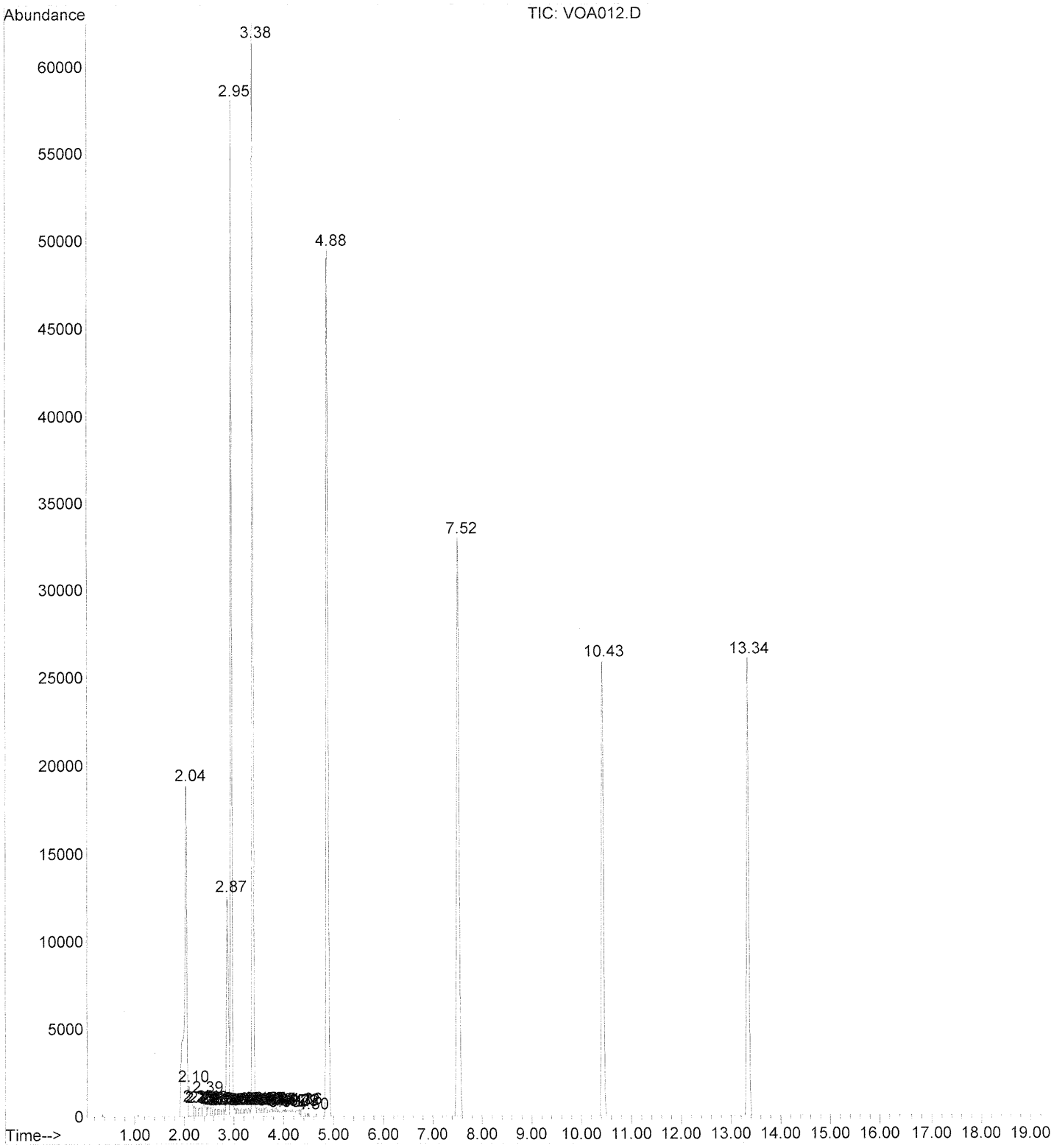
33 Carbon Tetrachloride	***** NOT FOUND *****			
34 2-Chloroethanol	***** NOT FOUND *****			
35 1,4-Difluorobenzene	3.38	100.0 ppb	ISTD2#	
36 1,2-Dichloropropane	***** NOT FOUND *****			
37 Trichloroethene	***** NOT FOUND *****			
38 Dibromomethane	***** NOT FOUND *****			
39 Bromodichloromethane	***** NOT FOUND *****			
40 Methyl methacrylate	***** NOT FOUND *****			
41 1,4-Dioxane	***** NOT FOUND *****			
42 cis-1,3-Dichloropropene	***** NOT FOUND *****			
43 2-Nitropropane	***** NOT FOUND *****			
44 trans-1,3-Dichloropropene	***** NOT FOUND *****			
45 Toluene	***** NOT FOUND *****			95 LR
46 Toluene-d8-surrogate	4.87	95.3 ppb		
47 1,1,2-Trichloroethane	***** NOT FOUND *****			
48 1,3-Dichloropropane	***** NOT FOUND *****			
49 Ethylmethacrylate	***** NOT FOUND *****			
50 Dibromochloromethane	***** NOT FOUND *****			
51 1,2-Dibromomethane	***** NOT FOUND *****			
52 Tetrachloroethene	***** NOT FOUND *****			
53 Chlorobenzene	***** NOT FOUND *****			
54 Chlorobenzene-d5	7.52	100.0 ppb	ISTD3	
55 1,1,1,2-Tetrachloroethane	***** NOT FOUND *****			
56 Ethylbenzene	***** NOT FOUND *****			
57 m/p-xylene	***** NOT FOUND *****			
58 Bromoform	***** NOT FOUND *****			
59 Styrene	***** NOT FOUND *****			
60 o-Xylene	***** NOT FOUND *****			
61 cis-1,4-Dichloro-2-butene	***** NOT FOUND *****			
62 trans-1,4-Dichloro-2-butene	***** NOT FOUND *****			
63 Bromofluorobenzene-surrogate	10.42	90.8 ppb		91 LR
64 1,1,2,2-Tetrachloroethane	***** NOT FOUND *****			
65 1,2,3-Trichloropropane	***** NOT FOUND *****			
66 Isopropylbenzene	***** NOT FOUND *****			
67 Bromobenzene	***** NOT FOUND *****			
68 2-Chlorotoluene	***** NOT FOUND *****			
69 n-Propylbenzene	***** NOT FOUND *****			
70 4-Chlorotoluene	***** NOT FOUND *****			
71 1,3,5-Trimethylbenzene	***** NOT FOUND *****			
72 Pentachloroethane	***** NOT FOUND *****			
73 tert-Butylbenzene	***** NOT FOUND *****			
74 1,2,4-Trimethylbenzene	***** NOT FOUND *****			
75 1,3-Dichlorobenzene	***** NOT FOUND *****			
76 1,4-Dichlorobenzene-d4	13.34	100.0 ppb	ISTD4	
77 sec-Butylbenzene	***** NOT FOUND *****			
78 1,4-Dichlorobenzene	***** NOT FOUND *****			
79 p-iso-Propyltoluene	***** NOT FOUND *****			
80 1,2-Dichlorobenzene	***** NOT FOUND *****			
81 n-Butylbenzene	***** NOT FOUND *****			
82 1,2-Bromo-3-chloropropane	***** NOT FOUND *****			
83 Nitobenzene	***** NOT FOUND *****			
84 1,2,4-Trichlorobenzene	***** NOT FOUND *****			
85 Naphthalene	***** NOT FOUND *****			
86 1,2,3-Trichlorobenzene	***** NOT FOUND *****			
87 Hexachlorobutadiene	***** NOT FOUND *****			

END OF REPORT

Qualifiers Not Satisfied

Wed May 24 16:01:02 2006

File : C:\HPCHEM\1\DATA\052406\VOA012.D
Operator : dmt
Acquired : 24 May 2006 4:08 pm using AcqMethod VOA0506
Instrument : GCD Syste
Sample Name: SB-19 6-10' H2O
Misc Info :
Vial Number: 1



Calibration Last Updated: Wed May 24 11:22:20 2006
Quantitation Report - Summary

Information from Data File:
File : C:\HPCHEM\1\DATA\052406\VOA012.D
Operator : dmt
Acquired : 24 May 2006 4:08 pm using AcqMethod VOA0506
Sample Name: SB-19 6-10' H2O
Misc Info :
Vial Number: 1
CurrentMeth: C:\HPCHEM\1\METHODS\VOA0506R.M

Quantitation Settings:
Non-Reference Peak Window: 10.00 Percent
Correlation Window: 0.02 Minutes
Default Multiplier: 1.00
Default Sample Concentration: 0.00
Peak Type Decoding: m -> Manually re-integrated

Number of Compounds in Database: 87

#	Compound	Ret Time	Concentration	Peak Type
1	Pentafluorobenzene	2.95	100.0 ppb	ISTD1
2	Dichlorofluoromethane	***** NOT FOUND *****		
3	Chloromethane	***** NOT FOUND *****		
4	Vinyl Chloride	***** NOT FOUND *****		
5	Bromomethane	***** NOT FOUND *****		
6	Chloroethane	***** NOT FOUND *****		
7	Trichlorofluoromethane	***** NOT FOUND *****		
8	Acetonitrile	***** NOT FOUND *****		
9	Diethyl ether	***** NOT FOUND *****		
10	1,1-Dichloroethene	***** NOT FOUND *****		
11	Acrylonitrile	***** NOT FOUND *****		
12	1,1,2-Trichlorotrifluoroethane	***** NOT FOUND *****		
13	Methylene Chloride	2.39	2.5 ppb	
14	Iodomethane	***** NOT FOUND *****		
15	trans-1,2-Dichloroethene	***** NOT FOUND *****		
16	MTBE	***** NOT FOUND *****		
17	Allyl chloride	***** NOT FOUND *****		
18	1,1-Dichloroethane	***** NOT FOUND *****		
19	Carbon disulfide	***** NOT FOUND *****		
20	Chloroprene	***** NOT FOUND *****		
21	Methacrylonitrile	***** NOT FOUND *****		
22	cis-1,2-Dichloroethene	***** NOT FOUND *****		
23	Methyl acrylate	***** NOT FOUND *****		
24	Chloroform	***** NOT FOUND *****		
25	2,2-Dichloropropane	***** NOT FOUND *****		
26	Dibromofluoromethane-surrogate	2.87	87.4 ppb	
27	Isobutyl alcohol	***** NOT FOUND *****		
28	Tetrahydrofuran	***** NOT FOUND *****		
29	1,1,1-Trichloroethane	***** NOT FOUND *****		
30	1,2-Dichloroethane	***** NOT FOUND *****		
31	1,1-Dichloropropene	***** NOT FOUND *****		
32	Benzene	***** NOT FOUND *****		

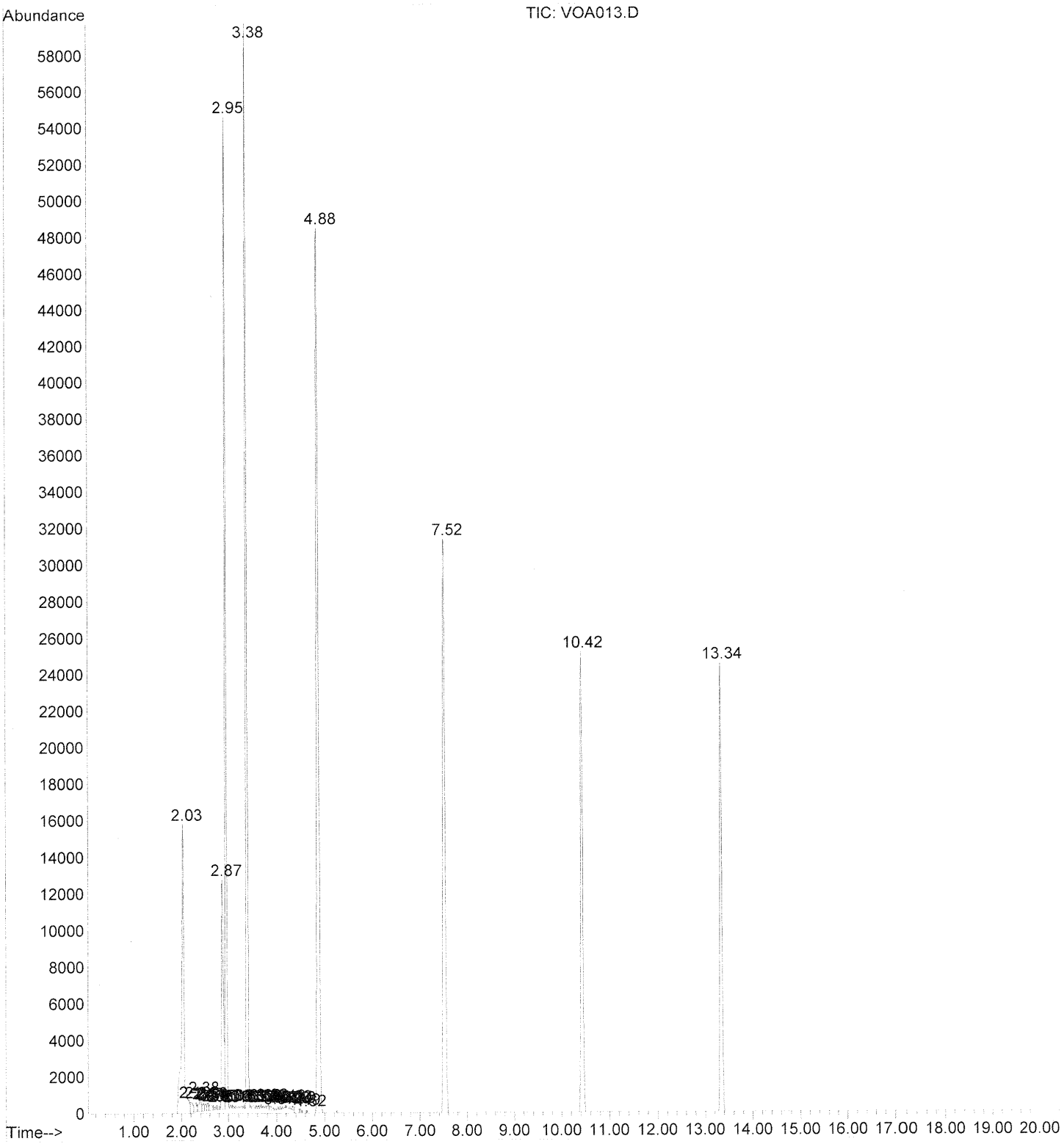
87 %R

33 Carbon Tetrachloride	***** NOT FOUND *****			
34 2-Chloroethanol	***** NOT FOUND *****			
35 1,4-Difluorobenzene	3.38	100.0 ppb	ISTD2	
36 1,2-Dichloropropane	***** NOT FOUND *****			
37 Trichloroethene	***** NOT FOUND *****			
38 Dibromomethane	***** NOT FOUND *****			
39 Bromodichloromethane	***** NOT FOUND *****			
40 Methyl methacrylate	***** NOT FOUND *****			
41 1,4-Dioxane	***** NOT FOUND *****			
42 cis-1,3-Dichloropropene	***** NOT FOUND *****			
43 2-Nitropropane	***** NOT FOUND *****			
44 trans-1,3-Dichloropropene	***** NOT FOUND *****			
45 Toluene	***** NOT FOUND *****			
46 Toluene-d8-surrogate	4.88	100.5 ppb		101 %R
47 1,1,2-Trichloroethane	***** NOT FOUND *****			
48 1,3-Dichloropropane	***** NOT FOUND *****			
49 Ethylmethacrylate	***** NOT FOUND *****			
50 Dibromochloromethane	***** NOT FOUND *****			
51 1,2-Dibromomethane	***** NOT FOUND *****			
52 Tetrachloroethene	***** NOT FOUND *****			
53 Chlorobenzene	***** NOT FOUND *****			
54 Chlorobenzene-d5	7.52	100.0 ppb	ISTD3	
55 1,1,1,2-Tetrachloroethane	***** NOT FOUND *****			
56 Ethylbenzene	***** NOT FOUND *****			
57 m/p-xylene	***** NOT FOUND *****			
58 Bromoform	***** NOT FOUND *****			
59 Styrene	***** NOT FOUND *****			
60 o-Xylene	***** NOT FOUND *****			
61 cis-1,4-Dichloro-2-butene	***** NOT FOUND *****			
62 trans-1,4-Dichloro-2-butene	***** NOT FOUND *****			
63 Bromofluorobenzene-surrogate	10.43	90.6 ppb		91 %R
64 1,1,2,2-Tetrachloroethane	***** NOT FOUND *****			
65 1,2,3-Trichloropropane	***** NOT FOUND *****			
66 Isopropylbenzene	***** NOT FOUND *****			
67 Bromobenzene	***** NOT FOUND *****			
68 2-Chlorotoluene	***** NOT FOUND *****			
69 n-Propylbenzene	***** NOT FOUND *****			
70 4-Chlorotoluene	***** NOT FOUND *****			
71 1,3,5-Trimethylbenzene	***** NOT FOUND *****			
72 Pentachloroethane	***** NOT FOUND *****			
73 tert-Butylbenzene	***** NOT FOUND *****			
74 1,2,4-Trimethylbenzene	***** NOT FOUND *****			
75 1,3-Dichlorobenzene	***** NOT FOUND *****			
76 1,4-Dichlorobenzene-d4	13.34	100.0 ppb	ISTD4	
77 sec-Butylbenzene	***** NOT FOUND *****			
78 1,4-Dichlorobenzene	***** NOT FOUND *****			
79 p-iso-Propyltoluene	***** NOT FOUND *****			
80 1,2-Dichlorobenzene	***** NOT FOUND *****			
81 n-Butylbenzene	***** NOT FOUND *****			
82 1,2-Bromo-3-chloropropane	***** NOT FOUND *****			
83 Nitobenzene	***** NOT FOUND *****			
84 1,2,4-Trichlorobenzene	***** NOT FOUND *****			
85 Naphthalene	***** NOT FOUND *****			
86 1,2,3-Trichlorobenzene	***** NOT FOUND *****			
87 Hexachlorobutadiene	***** NOT FOUND *****			

END OF REPORT

Wed May 24 16:29:23 2006

File : C:\HPCHEM\1\DATA\052406\VOA013.D
Operator : dmt
Acquired : 24 May 2006 4:35 pm using AcqMethod VOA0506
Instrument : GCD Syste
Sample Name: SB-20 10-14' H2O
Misc Info :
Vial Number: 1



Calibration Last Updated: Wed May 24 11:22:20 2006
Quantitation Report - Summary

Information from Data File:
File : C:\HPCHEM\1\DATA\052406\VOA013.D
Operator : dmt
Acquired : 24 May 2006 4:35 pm using AcqMethod VOA0506
Sample Name: SB-20 10-14' H2O
Misc Info :
Vial Number: 1
CurrentMeth: C:\HPCHEM\1\METHODS\VOA0506R.M

Quantitation Settings:
Non-Reference Peak Window: 10.00 Percent
Correlation Window: 0.02 Minutes
Default Multiplier: 1.00
Default Sample Concentration: 0.00
Peak Type Decoding: m -> Manually re-integrated

Number of Compounds in Database: 87

#	Compound	Ret Time	Concentration	Peak Type
1	Pentafluorobenzene	2.96	100.0 ppb	ISTD1
2	Dichlorofluoromethane	***** NOT FOUND *****		
3	Chloromethane	***** NOT FOUND *****		
4	Vinyl Chloride	***** NOT FOUND *****		
5	Bromomethane	***** NOT FOUND *****		
6	Chloroethane	***** NOT FOUND *****		
7	Trichlorofluoromethane	***** NOT FOUND *****		
8	Acetonitrile	***** NOT FOUND *****		
9	Diethyl ether	***** NOT FOUND *****		
10	1,1-Dichloroethene	***** NOT FOUND *****		
11	Acrylonitrile	***** NOT FOUND *****		
12	1,1,2-Trichlorotrifluoroethane	***** NOT FOUND *****		
13	Methylene Chloride	2.39	4.4 ppb	#
14	Iodomethane	***** NOT FOUND *****		
15	trans-1,2-Dichloroethene	***** NOT FOUND *****		
16	MTBE	***** NOT FOUND *****		
17	Allyl chloride	***** NOT FOUND *****		
18	1,1-Dichloroethane	***** NOT FOUND *****		
19	Carbon disulfide	***** NOT FOUND *****		
20	Chloroprene	***** NOT FOUND *****		
21	Methacrylonitrile	***** NOT FOUND *****		
22	cis-1,2-Dichloroethene	***** NOT FOUND *****		
23	Methyl acrylate	***** NOT FOUND *****		
24	Chloroform	***** NOT FOUND *****		
25	2,2-Dichloropropane	***** NOT FOUND *****		
26	Dibromofluoromethane-surrogate	2.87	95.1 ppb	
27	Isobutyl alcohol	***** NOT FOUND *****		
28	Tetrahydrofuran	***** NOT FOUND *****		
29	1,1,1-Trichloroethane	***** NOT FOUND *****		
30	1,2-Dichloroethane	***** NOT FOUND *****		
31	1,1-Dichloropropene	***** NOT FOUND *****		
32	Benzene	***** NOT FOUND *****		

95 %

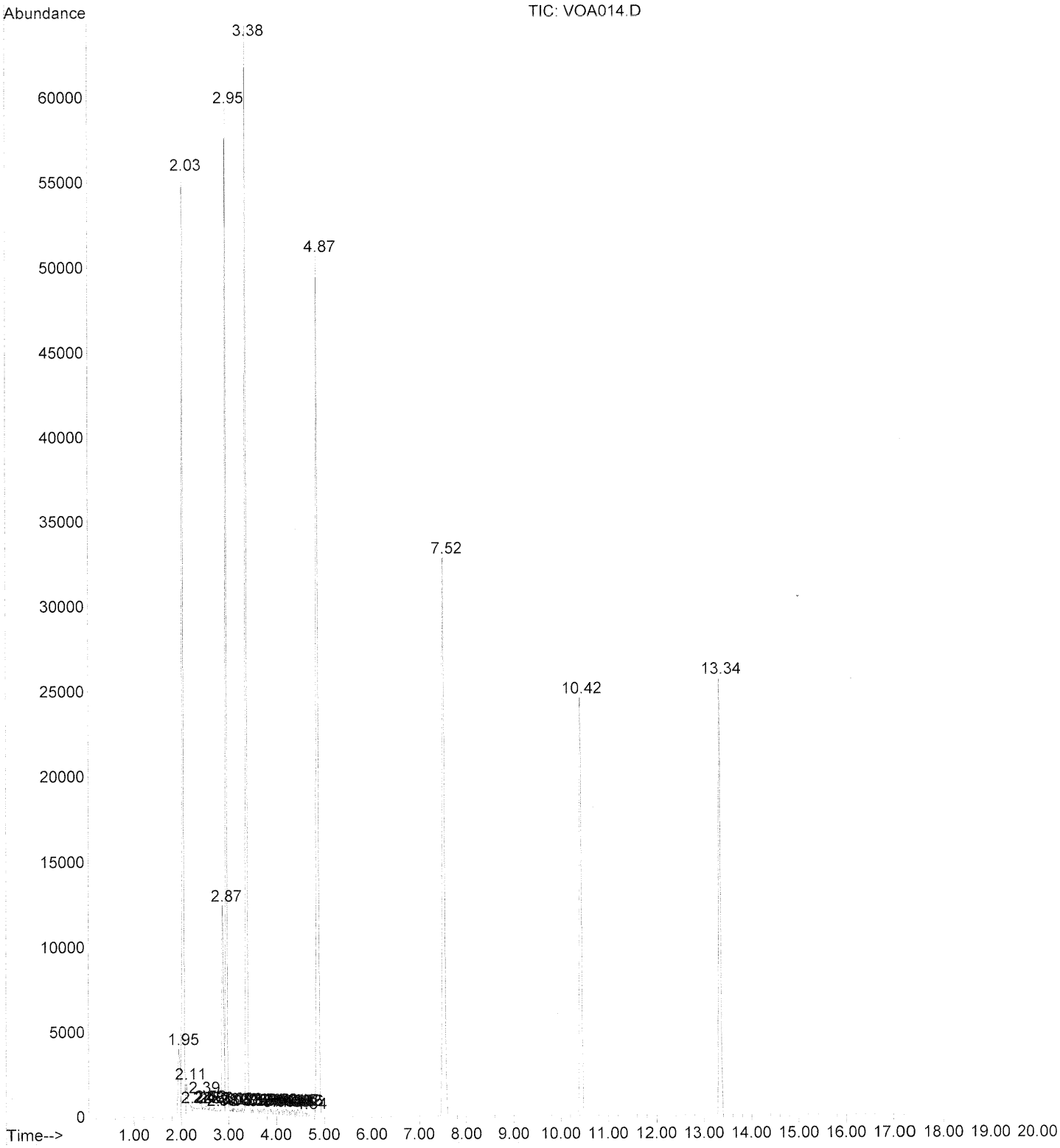
33 Carbon Tetrachloride	***** NOT FOUND *****		
34 2-Chloroethanol	***** NOT FOUND *****		
35 1,4-Difluorobenzene	3.38	100.0 ppb	ISTD2
36 1,2-Dichloropropane	***** NOT FOUND *****		
37 Trichloroethene	***** NOT FOUND *****		
38 Dibromomethane	***** NOT FOUND *****		
39 Bromodichloromethane	***** NOT FOUND *****		
40 Methyl methacrylate	***** NOT FOUND *****		
41 1,4-Dioxane	***** NOT FOUND *****		
42 cis-1,3-Dichloropropene	***** NOT FOUND *****		
43 2-Nitropropane	***** NOT FOUND *****		
44 trans-1,3-Dichloropropene	***** NOT FOUND *****		
45 Toluene	***** NOT FOUND *****		
46 Toluene-d8-surrogate	4.88	104.3 ppb	104 %
47 1,1,2-Trichloroethane	***** NOT FOUND *****		
48 1,3-Dichloropropane	***** NOT FOUND *****		
49 Ethylmethacrylate	***** NOT FOUND *****		
50 Dibromochloromethane	***** NOT FOUND *****		
51 1,2-Dibromomethane	***** NOT FOUND *****		
52 Tetrachloroethene	***** NOT FOUND *****		
53 Chlorobenzene	***** NOT FOUND *****		
54 Chlorobenzene-d5	7.53	100.0 ppb	ISTD3
55 1,1,1,2-Tetrachloroethane	***** NOT FOUND *****		
56 Ethylbenzene	***** NOT FOUND *****		
57 m/p-xylene	***** NOT FOUND *****		
58 Bromoform	***** NOT FOUND *****		
59 Styrene	***** NOT FOUND *****		
60 o-Xylene	***** NOT FOUND *****		
61 cis-1,4-Dichloro-2-butene	***** NOT FOUND *****		
62 trans-1,4-Dichloro-2-butene	***** NOT FOUND *****		
63 Bromofluorobenzene-surrogate	10.42	92.5 ppb	93 %
64 1,1,2,2-Tetrachloroethane	***** NOT FOUND *****		
65 1,2,3-Trichloropropane	***** NOT FOUND *****		
66 Isopropylbenzene	***** NOT FOUND *****		
67 Bromobenzene	***** NOT FOUND *****		
68 2-Chlorotoluene	***** NOT FOUND *****		
69 n-Propylbenzene	***** NOT FOUND *****		
70 4-Chlorotoluene	***** NOT FOUND *****		
71 1,3,5-Trimethylbenzene	***** NOT FOUND *****		
72 Pentachloroethane	***** NOT FOUND *****		
73 tert-Butylbenzene	***** NOT FOUND *****		
74 1,2,4-Trimethylbenzene	***** NOT FOUND *****		
75 1,3-Dichlorobenzene	***** NOT FOUND *****		
76 1,4-Dichlorobenzene-d4	13.33	100.0 ppb	ISTD4
77 sec-Butylbenzene	***** NOT FOUND *****		
78 1,4-Dichlorobenzene	***** NOT FOUND *****		
79 p-iso-Propyltoluene	***** NOT FOUND *****		
80 1,2-Dichlorobenzene	***** NOT FOUND *****		
81 n-Butylbenzene	***** NOT FOUND *****		
82 1,2-Bromo-3-chloropropane	***** NOT FOUND *****		
83 Nitobenzene	***** NOT FOUND *****		
84 1,2,4-Trichlorobenzene	***** NOT FOUND *****		
85 Naphthalene	***** NOT FOUND *****		
86 1,2,3-Trichlorobenzene	***** NOT FOUND *****		
87 Hexachlorobutadiene	***** NOT FOUND *****		

END OF REPORT

Qualifiers Not Satisfied

Wed May 24 16:56:47 2006

File : C:\HPCHEM\1\DATA\052406\VOA014.D
Operator : dmt
Acquired : 24 May 2006 5:05 pm using AcqMethod VOA0506
Instrument : GCD Syste
Sample Name: SB-18 10-14' H2O
Misc Info :
Vial Number: 1



Calibration Last Updated: Wed May 24 11:22:20 2006
Quantitation Report - Summary

Information from Data File:
File : C:\HPCHEM\1\DATA\052406\VOA014.D
Operator : dmt
Acquired : 24 May 2006 5:05 pm using AcqMethod VOA0506
Sample Name: SB-18 10-14' H2O
Misc Info :
Vial Number: 1
CurrentMeth: C:\HPCHEM\1\METHODS\VOA0506R.M

Quantitation Settings:
Non-Reference Peak Window: 10.00 Percent
Correlation Window: 0.02 Minutes
Default Multiplier: 1.00
Default Sample Concentration: 0.00
Peak Type Decoding: m -> Manually re-integrated

Number of Compounds in Database: 87

#	Compound	Ret Time	Concentration	Peak Type
1	Pentafluorobenzene	2.95	100.0 ppb	ISTD1
2	Dichlorofluoromethane	***** NOT FOUND *****		
3	Chloromethane	***** NOT FOUND *****		
4	Vinyl Chloride	***** NOT FOUND *****		
5	Bromomethane	***** NOT FOUND *****		
6	Chloroethane	***** NOT FOUND *****		
7	Trichlorofluoromethane	***** NOT FOUND *****		
8	Acetonitrile	***** NOT FOUND *****		
9	Diethyl ether	***** NOT FOUND *****		
10	1,1-Dichloroethene	***** NOT FOUND *****		
11	Acrylonitrile	***** NOT FOUND *****		
12	1,1,2-Trichlorotrifluoroethane	***** NOT FOUND *****		
13	Methylene Chloride	2.38	3.8 ppb	#
14	Iodomethane	***** NOT FOUND *****		
15	trans-1,2-Dichloroethene	***** NOT FOUND *****		
16	MTBE	***** NOT FOUND *****		
17	Allyl chloride	***** NOT FOUND *****		
18	1,1-Dichloroethane	***** NOT FOUND *****		
19	Carbon disulfide	***** NOT FOUND *****		
20	Chloroprene	***** NOT FOUND *****		
21	Methacrylonitrile	***** NOT FOUND *****		
22	cis-1,2-Dichloroethene	***** NOT FOUND *****		
23	Methyl acrylate	***** NOT FOUND *****		
24	Chloroform	***** NOT FOUND *****		
25	2,2-Dichloropropane	***** NOT FOUND *****		
26	Dibromofluoromethane-surrogate	2.87	87.0 ppb	
27	Isobutyl alcohol	***** NOT FOUND *****		
28	Tetrahydrofuran	***** NOT FOUND *****		
29	1,1,1-Trichloroethane	***** NOT FOUND *****		
30	1,2-Dichloroethane	***** NOT FOUND *****		
31	1,1-Dichloropropene	***** NOT FOUND *****		
32	Benzene	***** NOT FOUND *****		

87 %R

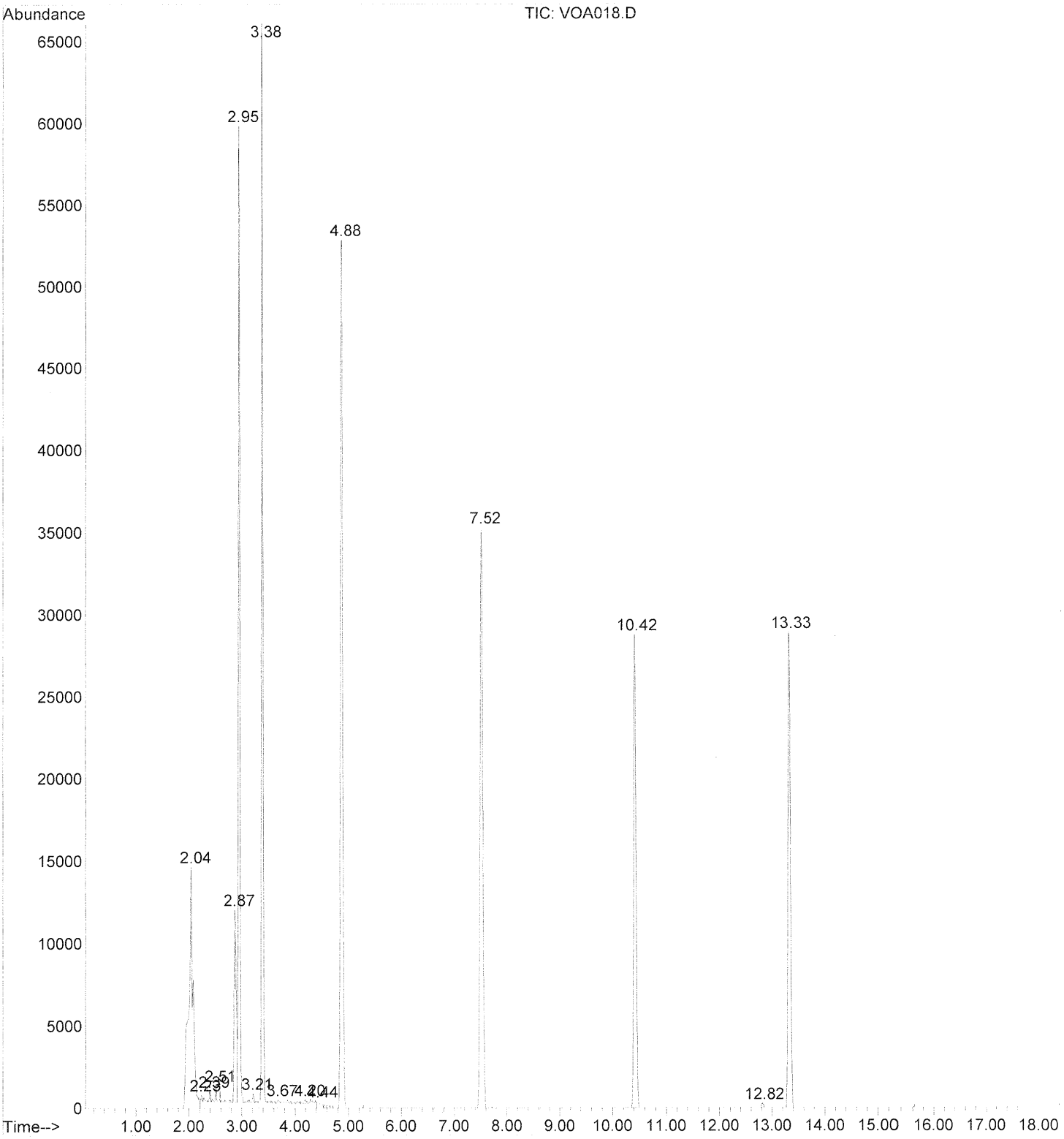
33 Carbon Tetrachloride	***** NOT FOUND *****			
34 2-Chloroethanol	***** NOT FOUND *****			
35 1,4-Difluorobenzene	3.38	100.0 ppb	ISTD2	
36 1,2-Dichloropropane	***** NOT FOUND *****			
37 Trichloroethene	***** NOT FOUND *****			
38 Dibromomethane	***** NOT FOUND *****			
39 Bromodichloromethane	***** NOT FOUND *****			
40 Methyl methacrylate	***** NOT FOUND *****			
41 1,4-Dioxane	***** NOT FOUND *****			
42 cis-1,3-Dichloropropene	***** NOT FOUND *****			
43 2-Nitropropane	***** NOT FOUND *****			
44 trans-1,3-Dichloropropene	***** NOT FOUND *****			
45 Toluene	***** NOT FOUND *****			
46 Toluene-d8-surrogate	4.87	99.5 ppb		100 %R
47 1,1,2-Trichloroethane	***** NOT FOUND *****			
48 1,3-Dichloropropane	***** NOT FOUND *****			
49 Ethylmethacrylate	***** NOT FOUND *****			
50 Dibromochloromethane	***** NOT FOUND *****			
51 1,2-Dibromomethane	***** NOT FOUND *****			
52 Tetrachloroethene	***** NOT FOUND *****			
53 Chlorobenzene	***** NOT FOUND *****			
54 Chlorobenzene-d5	7.51	100.0 ppb	ISTD3	
55 1,1,1,2-Tetrachloroethane	***** NOT FOUND *****			
56 Ethylbenzene	***** NOT FOUND *****			
57 m/p-xylene	***** NOT FOUND *****			
58 Bromoform	***** NOT FOUND *****			
59 Styrene	***** NOT FOUND *****			
60 o-Xylene	***** NOT FOUND *****			
61 cis-1,4-Dichloro-2-butene	***** NOT FOUND *****			
62 trans-1,4-Dichloro-2-butene	***** NOT FOUND *****			
63 Bromofluorobenzene-surrogate	10.42	85.4 ppb		85 %R
64 1,1,2,2-Tetrachloroethane	***** NOT FOUND *****			
65 1,2,3-Trichloropropane	***** NOT FOUND *****			
66 Isopropylbenzene	***** NOT FOUND *****			
67 Bromobenzene	***** NOT FOUND *****			
68 2-Chlorotoluene	***** NOT FOUND *****			
69 n-Propylbenzene	***** NOT FOUND *****			
70 4-Chlorotoluene	***** NOT FOUND *****			
71 1,3,5-Trimethylbenzene	***** NOT FOUND *****			
72 Pentachloroethane	***** NOT FOUND *****			
73 tert-Butylbenzene	***** NOT FOUND *****			
74 1,2,4-Trimethylbenzene	***** NOT FOUND *****			
75 1,3-Dichlorobenzene	***** NOT FOUND *****			
76 1,4-Dichlorobenzene-d4	13.34	100.0 ppb	ISTD4	
77 sec-Butylbenzene	***** NOT FOUND *****			
78 1,4-Dichlorobenzene	***** NOT FOUND *****			
79 p-iso-Propyltoluene	***** NOT FOUND *****			
80 1,2-Dichlorobenzene	***** NOT FOUND *****			
81 n-Butylbenzene	***** NOT FOUND *****			
82 1,2-Bromo-3-chloropropane	***** NOT FOUND *****			
83 Nitobenzene	***** NOT FOUND *****			
84 1,2,4-Trichlorobenzene	***** NOT FOUND *****			
85 Naphthalene	***** NOT FOUND *****			
86 1,2,3-Trichlorobenzene	***** NOT FOUND *****			
87 Hexachlorobutadiene	***** NOT FOUND *****			

END OF REPORT

Qualifiers Not Satisfied

Thu May 25 07:34:47 2006

File : C:\HPCHEM\1\DATA\052406\VOA018.D
Operator : dmt
Acquired : 25 May 2006 9:22 am using AcqMethod VOA0506
Instrument : GCD Syste
Sample Name: SB-32 6-10' H2O
Misc Info :
Vial Number: 1



Calibration Last Updated: Thu May 25 08:53:45 2006
Quantitation Report - Summary

Information from Data File:
File : C:\HPCHEM\1\DATA\052406\VOA018.D
Operator : dmt
Acquired : 25 May 2006 9:22 am using AcqMethod VOA0506
Sample Name: SB-32 6-10' H2O
Misc Info :
Vial Number: 1
CurrentMeth: C:\HPCHEM\1\METHODS\VOA0506R.M

Quantitation Settings:
Non-Reference Peak Window: 10.00 Percent
Correlation Window: 0.02 Minutes
Default Multiplier: 1.00
Default Sample Concentration: 0.00
Peak Type Decoding: m -> Manually re-integrated

Number of Compounds in Database: 87

#	Compound	Ret Time	Concentration	Peak Type
1	Pentafluorobenzene	2.95	100.0 ppb	ISTD1
2	Dichlorofluoromethane	~~2.09~~	~~2.4 ppb~~	~~#~~
3	Chloromethane	***** NOT FOUND *****		
4	Vinyl Chloride	***** NOT FOUND *****		
5	Bromomethane	***** NOT FOUND *****		
6	Chloroethane	2.08	4.8 ppb	#
7	Trichlorofluoromethane	***** NOT FOUND *****		
8	Acetonitrile	***** NOT FOUND *****		
9	Diethyl ether	***** NOT FOUND *****		
10	1,1-Dichloroethene	***** NOT FOUND *****		
11	Acrylonitrile	***** NOT FOUND *****		
12	1,1,2-Trichlorotrifluoroethane	***** NOT FOUND *****		
13	Methylene Chloride	2.39	4.2 ppb	#
14	Iodomethane	***** NOT FOUND *****		
15	trans-1,2-Dichloroethene	***** NOT FOUND *****		
16	MTBE	***** NOT FOUND *****		
17	Allyl chloride	***** NOT FOUND *****		
18	1,1-Dichloroethane	***** NOT FOUND *****		
19	Carbon disulfide	***** NOT FOUND *****		
20	Chloroprene	***** NOT FOUND *****		
21	Methacrylonitrile	***** NOT FOUND *****		
22	cis-1,2-Dichloroethene	***** NOT FOUND *****		
23	Methyl acrylate	***** NOT FOUND *****		
24	Chloroform	***** NOT FOUND *****		
25	2,2-Dichloropropane	***** NOT FOUND *****		
26	Dibromofluoromethane-surrogate	2.87	91.2 ppb	
27	Isobutyl alcohol	***** NOT FOUND *****		
28	Tetrahydrofuran	***** NOT FOUND *****		
29	1,1,1-Trichloroethane	***** NOT FOUND *****		
30	1,2-Dichloroethane	***** NOT FOUND *****		
31	1,1-Dichloropropene	***** NOT FOUND *****		
32	Benzene	3.22	2.1 ppb	#

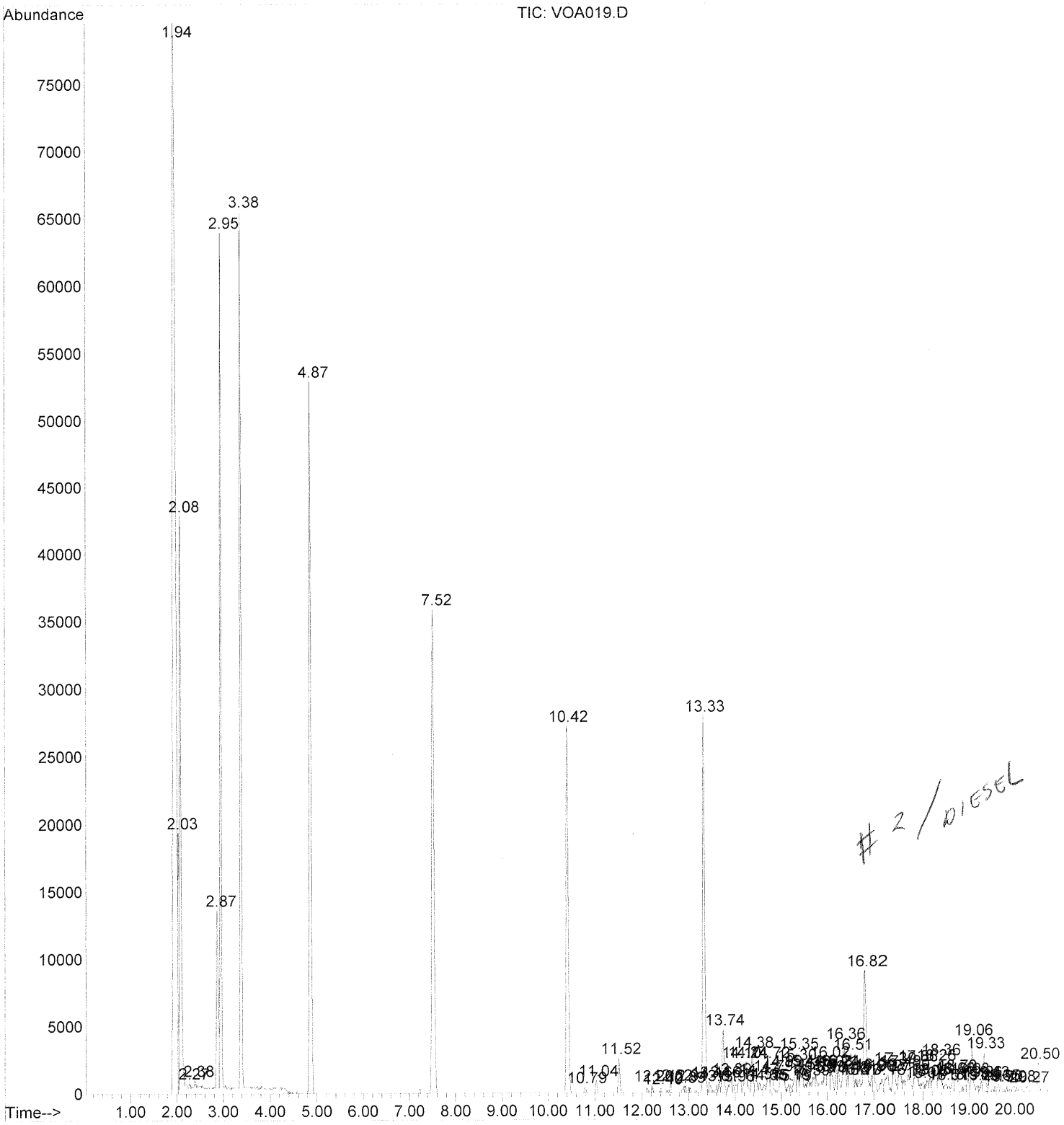
33 Carbon Tetrachloride	***** NOT FOUND *****			
34 2-Chloroethanol	***** NOT FOUND *****			
35 1,4-Difluorobenzene	3.38	100.0 ppb		ISTD2
36 1,2-Dichloropropane	***** NOT FOUND *****			
37 Trichloroethene	***** NOT FOUND *****			
38 Dibromomethane	***** NOT FOUND *****			
39 Bromodichloromethane	***** NOT FOUND *****			
40 Methyl methacrylate	***** NOT FOUND *****			
41 1,4-Dioxane	***** NOT FOUND *****			
42 cis-1,3-Dichloropropene	***** NOT FOUND *****			
43 2-Nitropropane	***** NOT FOUND *****			
44 trans-1,3-Dichloropropene	***** NOT FOUND *****			
45 Toluene	***** NOT FOUND *****			
46 Toluene-d8-surrogate	4.88	106.6 ppb		107%
47 1,1,2-Trichloroethane	***** NOT FOUND *****			
48 1,3-Dichloropropane	***** NOT FOUND *****			
49 Ethylmethacrylate	***** NOT FOUND *****			
50 Dibromochloromethane	***** NOT FOUND *****			
51 1,2-Dibromomethane	***** NOT FOUND *****			
52 Tetrachloroethene	***** NOT FOUND *****			
53 Chlorobenzene	***** NOT FOUND *****			
54 Chlorobenzene-d5	7.53	100.0 ppb		ISTD3
55 1,1,1,2-Tetrachloroethane	***** NOT FOUND *****			
56 Ethylbenzene	***** NOT FOUND *****			
57 m/p-xylene	***** NOT FOUND *****			
58 Bromoform	***** NOT FOUND *****			
59 Styrene	***** NOT FOUND *****			
60 o-Xylene	***** NOT FOUND *****			
61 cis-1,4-Dichloro-2-butene	***** NOT FOUND *****			
62 trans-1,4-Dichloro-2-butene	***** NOT FOUND *****			
63 Bromofluorobenzene-surrogate	10.42	94.7 ppb		95%
64 1,1,2,2-Tetrachloroethane	***** NOT FOUND *****			
65 1,2,3-Trichloropropane	***** NOT FOUND *****			
66 Isopropylbenzene	***** NOT FOUND *****			
67 Bromobenzene	***** NOT FOUND *****			
68 2-Chlorotoluene	***** NOT FOUND *****			
69 n-Propylbenzene	***** NOT FOUND *****			
70 4-Chlorotoluene	***** NOT FOUND *****			
71 1,3,5-Trimethylbenzene	***** NOT FOUND *****			
72 Pentachloroethane	***** NOT FOUND *****			
73 tert-Butylbenzene	***** NOT FOUND *****			
74 1,2,4-Trimethylbenzene	12.82	2.9 ppb	#	2.9
75 1,3-Dichlorobenzene	***** NOT FOUND *****			
76 1,4-Dichlorobenzene-d4	13.33	100.0 ppb		ISTD4
77 sec-Butylbenzene	***** NOT FOUND *****			
78 1,4-Dichlorobenzene	***** NOT FOUND *****			
79 p-iso-Propyltoluene	***** NOT FOUND *****			
80 1,2-Dichlorobenzene	***** NOT FOUND *****			
81 n-Butylbenzene	***** NOT FOUND *****			
82 1,2-Bromo-3-chloropropane	***** NOT FOUND *****			
83 Nitobenzene	***** NOT FOUND *****			
84 1,2,4-Trichlorobenzene	***** NOT FOUND *****			
85 Naphthalene	***** NOT FOUND *****			
86 1,2,3-Trichlorobenzene	***** NOT FOUND *****			
87 Hexachlorobutadiene	***** NOT FOUND *****			

END OF REPORT

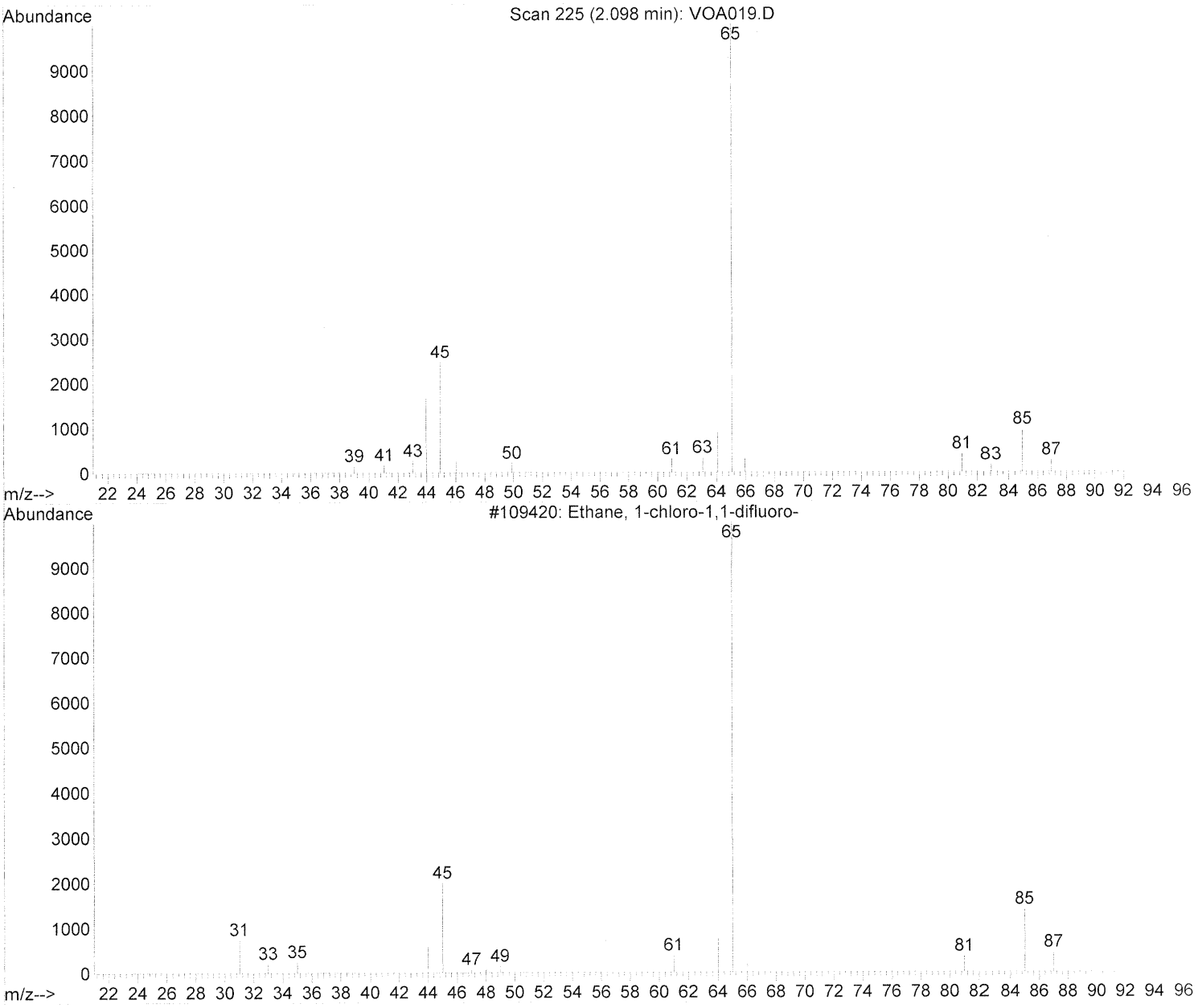
Qualifiers Not Satisfied

Thu May 25 09:43:48 2006

File : C:\HPCHEM\1\DATA\052406\VOA019.D
Operator : dmt
Acquired : 25 May 2006 9:49 am using AcqMethod VOA0506
Instrument : GCD Syste
Sample Name: SB-33 6-10' H2O
Misc Info :
Vial Number: 1



Library Searched : C:\DATABASE\NIST129K.L
Quality : 83
ID : Ethane, 1-chloro-1,1-difluoro-



Calibration Last Updated: Thu May 25 08:53:45 2006
Quantitation Report - Summary

Information from Data File:
File : C:\HPCHEM\1\DATA\052406\VOA019.D
Operator : dmt
Acquired : 25 May 2006 9:49 am using AcqMethod VOA0506
Sample Name: SB-33 6-10' H2O
Misc Info :
Vial Number: 1
CurrentMeth: C:\HPCHEM\1\METHODS\VOA0506R.M

Quantitation Settings:
Non-Reference Peak Window: 10.00 Percent
Correlation Window: 0.02 Minutes
Default Multiplier: 1.00
Default Sample Concentration: 0.00
Peak Type Decoding: m -> Manually re-integrated

Number of Compounds in Database: 87

#	Compound	Ret Time	Concentration	Peak Type
1	Pentafluorobenzene	2.95	100.0 ppb	ISTD1
2	Dichlorofluoromethane	2.08	13.1 ppb	
3	Chloromethane	~~2.08~~	~~1.4 ppb~~ #	
4	Vinyl Chloride	***** NOT FOUND *****		
5	Bromomethane	***** NOT FOUND *****		
6	Chloroethane	~~2.07~~	~~28.1 ppb~~	
7	Trichlorofluoromethane	***** NOT FOUND *****		
8	Acetonitrile	***** NOT FOUND *****		
9	Diethyl ether	***** NOT FOUND *****		
10	1,1-Dichloroethene	***** NOT FOUND *****		
11	Acrylonitrile	***** NOT FOUND *****		
12	1,1,2-Trichlorotrifluoroethane	***** NOT FOUND *****		
13	Methylene Chloride	2.38	3.8 ppb #	
14	Iodomethane	***** NOT FOUND *****		
15	trans-1,2-Dichloroethene	***** NOT FOUND *****		
16	MTBE	***** NOT FOUND *****		
17	Allyl chloride	***** NOT FOUND *****		
18	1,1-Dichloroethane	***** NOT FOUND *****		
19	Carbon disulfide	***** NOT FOUND *****		
20	Chloroprene	***** NOT FOUND *****		
21	Methacrylonitrile	***** NOT FOUND *****		
22	cis-1,2-Dichloroethene	***** NOT FOUND *****		
23	Methyl acrylate	***** NOT FOUND *****		
24	Chloroform	***** NOT FOUND *****		
25	2,2-Dichloropropane	***** NOT FOUND *****		
26	Dibromofluoromethane-surrogate	2.87	93.9 ppb	
27	Isobutyl alcohol	***** NOT FOUND *****		
28	Tetrahydrofuran	***** NOT FOUND *****		
29	1,1,1-Trichloroethane	***** NOT FOUND *****		
30	1,2-Dichloroethane	***** NOT FOUND *****		
31	1,1-Dichloropropene	***** NOT FOUND *****		
32	Benzene	***** NOT FOUND *****		

13 (CHLORO-DIFLUORO METHANE)

#2/DIESEL

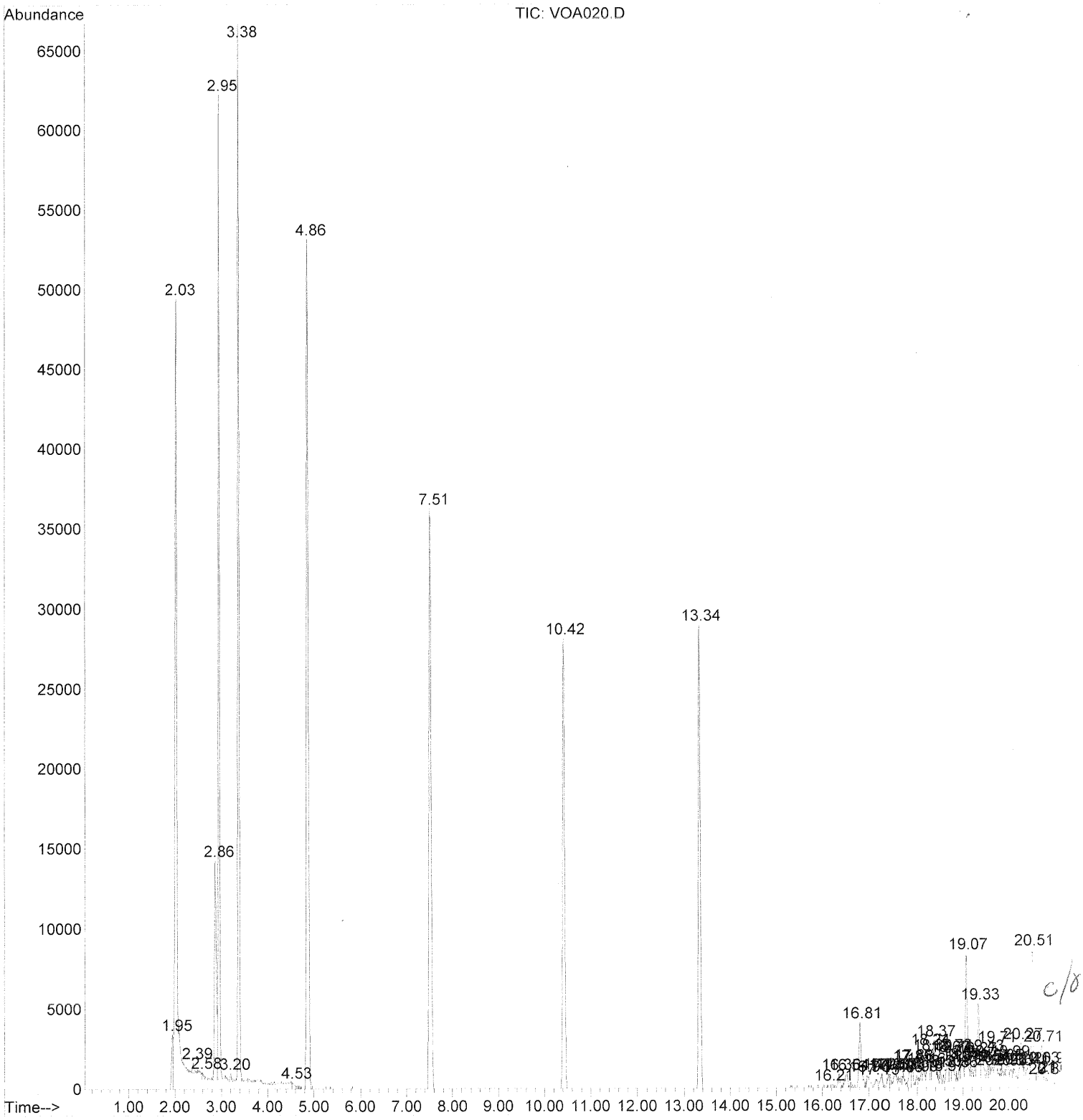
33 Carbon Tetrachloride	***** NOT FOUND *****		
34 2-Chloroethanol	***** NOT FOUND *****		
35 1,4-Difluorobenzene	3.38	100.0 ppb	ISTD2
36 1,2-Dichloropropane	***** NOT FOUND *****		
37 Trichloroethene	***** NOT FOUND *****		
38 Dibromomethane	***** NOT FOUND *****		
39 Bromodichloromethane	***** NOT FOUND *****		
40 Methyl methacrylate	***** NOT FOUND *****		
41 1,4-Dioxane	***** NOT FOUND *****		
42 cis-1,3-Dichloropropene	***** NOT FOUND *****		
43 2-Nitropropane	***** NOT FOUND *****		
44 trans-1,3-Dichloropropene	***** NOT FOUND *****		
45 Toluene	***** NOT FOUND *****		
46 Toluene-d8-surrogate	4.87	102.6 ppb	103%
47 1,1,2-Trichloroethane	***** NOT FOUND *****		
48 1,3-Dichloropropane	***** NOT FOUND *****		
49 Ethylmethacrylate	***** NOT FOUND *****		
50 Dibromochloromethane	***** NOT FOUND *****		
51 1,2-Dibromomethane	***** NOT FOUND *****		
52 Tetrachloroethene	***** NOT FOUND *****		
53 Chlorobenzene	***** NOT FOUND *****		
54 Chlorobenzene-d5	7.52	100.0 ppb	ISTD3
55 1,1,1,2-Tetrachloroethane	***** NOT FOUND *****		
56 Ethylbenzene	***** NOT FOUND *****		
57 m/p-xylene	***** NOT FOUND *****		
58 Bromoform	***** NOT FOUND *****		
59 Styrene	***** NOT FOUND *****		
60 o-Xylene	***** NOT FOUND *****		
61 cis-1,4-Dichloro-2-butene	***** NOT FOUND *****		
62 trans-1,4-Dichloro-2-butene	***** NOT FOUND *****		
63 Bromofluorobenzene-surrogate	10.42	91.7 ppb	92%
64 1,1,2,2-Tetrachloroethane	***** NOT FOUND *****		
65 1,2,3-Trichloropropane	***** NOT FOUND *****		
66 Isopropylbenzene	***** NOT FOUND *****		
67 Bromobenzene	***** NOT FOUND *****		
68 2-Chlorotoluene	***** NOT FOUND *****		
69 n-Propylbenzene	***** NOT FOUND *****		
70 4-Chlorotoluene	***** NOT FOUND *****		
71 1,3,5-Trimethylbenzene	***** NOT FOUND *****		
72 Pentachloroethane	***** NOT FOUND *****		
73 tert-Butylbenzene	***** NOT FOUND *****		
74 1,2,4-Trimethylbenzene	***** NOT FOUND *****		
75 1,3-Dichlorobenzene	***** NOT FOUND *****		
76 1,4-Dichlorobenzene-d4	13.34	100.0 ppb	ISTD4
77 sec-Butylbenzene	***** NOT FOUND *****		
78 1,4-Dichlorobenzene	***** NOT FOUND *****		
79 p-iso-Propyltoluene	***** NOT FOUND *****		
80 1,2-Dichlorobenzene	***** NOT FOUND *****		
81 n-Butylbenzene	***** NOT FOUND *****		
82 1,2-Bromo-3-chloropropane	***** NOT FOUND *****		
83 Nitobenzene	***** NOT FOUND *****		
84 1,2,4-Trichlorobenzene	***** NOT FOUND *****		
85 Naphthalene	***** NOT FOUND *****		
86 1,2,3-Trichlorobenzene	***** NOT FOUND *****		
87 Hexachlorobutadiene	***** NOT FOUND *****		

END OF REPORT

Qualifiers Not Satisfied

Thu May 25 10:13:00 2006

File : C:\HPCHEM\1\DATA\052406\VOA020.D
Operator : dmt
Acquired : 25 May 2006 10:18 am using AcqMethod VOA0506
Instrument : GCD Syste
Sample Name: SB-1 6-10' H2O
Misc Info :
Vial Number: 1



Calibration Last Updated: Thu May 25 08:53:45 2006
Quantitation Report - Summary

Information from Data File:
File : C:\HPCHEM\1\DATA\052406\VOA020.D
Operator : dmt
Acquired : 25 May 2006 10:18 am using AcqMethod VOA0506
Sample Name: SB-1 6-10' H2O
Misc Info :
Vial Number: 1
CurrentMeth: C:\HPCHEM\1\METHODS\VOA0506R.M

Quantitation Settings:
Non-Reference Peak Window: 10.00 Percent
Correlation Window: 0.02 Minutes
Default Multiplier: 1.00
Default Sample Concentration: 0.00
Peak Type Decoding: m -> Manually re-integrated

Number of Compounds in Database: 87

#	Compound	Ret Time	Concentration	Peak Type
1	Pentafluorobenzene	2.95	100.0 ppb	ISTD1
2	Dichlorofluoromethane	***** NOT FOUND *****		
3	Chloromethane	***** NOT FOUND *****		
4	Vinyl Chloride	***** NOT FOUND *****		
5	Bromomethane	***** NOT FOUND *****		
6	Chloroethane	***** NOT FOUND *****		
7	Trichlorofluoromethane	***** NOT FOUND *****		
8	Acetonitrile	***** NOT FOUND *****		
9	Diethyl ether	***** NOT FOUND *****		
10	1,1-Dichloroethene	***** NOT FOUND *****		
11	Acrylonitrile	***** NOT FOUND *****		
12	1,1,2-Trichlorotrifluoroethane	***** NOT FOUND *****		
13	Methylene Chloride	2.39	4.7 ppb	#
14	Iodomethane	***** NOT FOUND *****		
15	trans-1,2-Dichloroethene	***** NOT FOUND *****		
16	MTBE	***** NOT FOUND *****		
17	Allyl chloride	***** NOT FOUND *****		
18	1,1-Dichloroethane	***** NOT FOUND *****		
19	Carbon disulfide	***** NOT FOUND *****		
20	Chloroprene	***** NOT FOUND *****		
21	Methacrylonitrile	***** NOT FOUND *****		
22	cis-1,2-Dichloroethene	***** NOT FOUND *****		
23	Methyl acrylate	***** NOT FOUND *****		
24	Chloroform	***** NOT FOUND *****		
25	2,2-Dichloropropane	***** NOT FOUND *****		
26	Dibromofluoromethane-surrogate	2.86	97.0 ppb	
27	Isobutyl alcohol	***** NOT FOUND *****		
28	Tetrahydrofuran	***** NOT FOUND *****		
29	1,1,1-Trichloroethane	***** NOT FOUND *****		
30	1,2-Dichloroethane	***** NOT FOUND *****		
31	1,1-Dichloropropene	***** NOT FOUND *****		
32	Benzene	***** NOT FOUND *****		

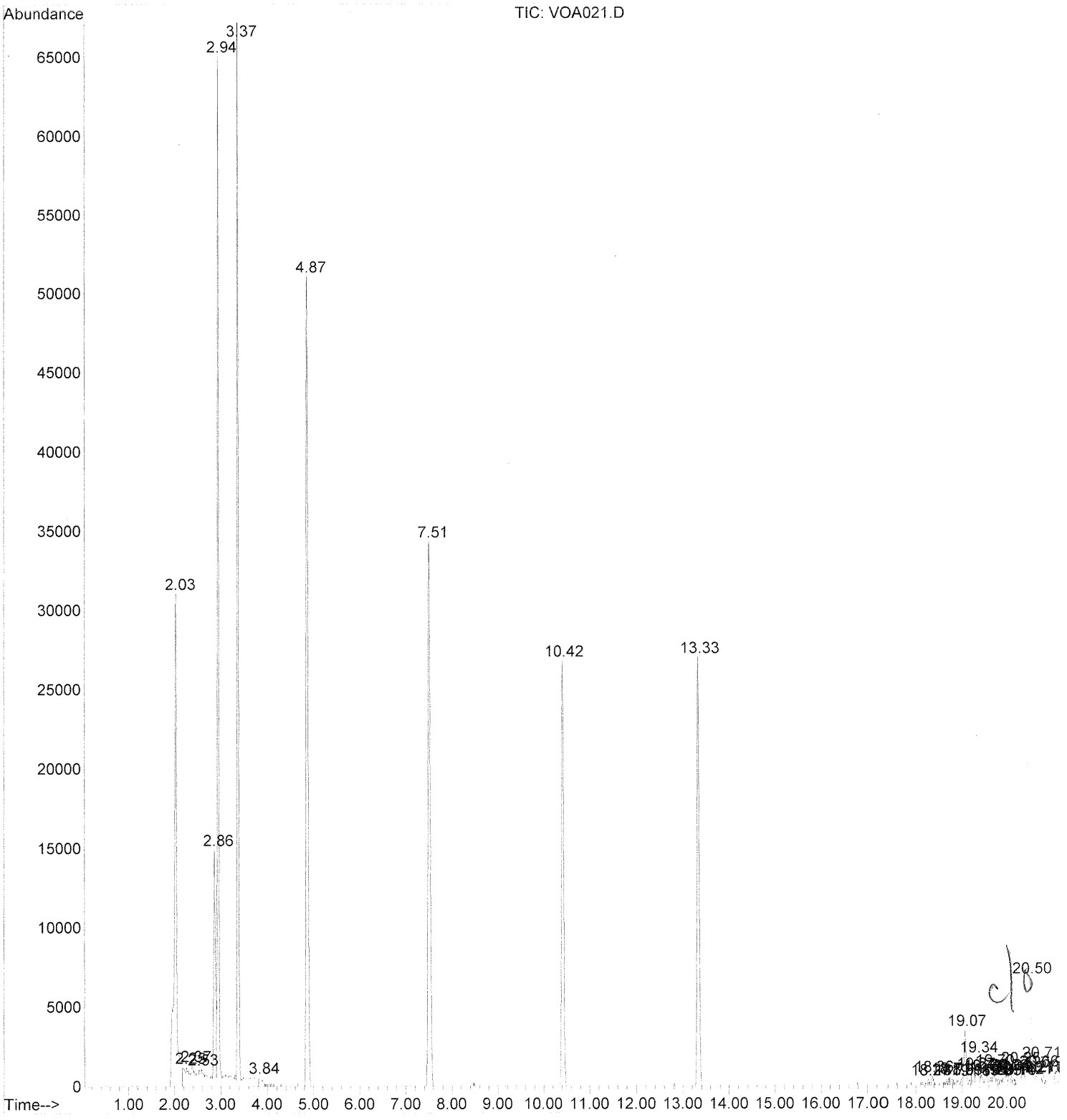
33 Carbon Tetrachloride	***** NOT FOUND *****		
34 2-Chloroethanol	***** NOT FOUND *****		
35 1,4-Difluorobenzene	3.38	100.0 ppb	ISTD2
36 1,2-Dichloropropane	***** NOT FOUND *****		
37 Trichloroethene	***** NOT FOUND *****		
38 Dibromomethane	***** NOT FOUND *****		
39 Bromodichloromethane	***** NOT FOUND *****		
40 Methyl methacrylate	***** NOT FOUND *****		
41 1,4-Dioxane	***** NOT FOUND *****		
42 cis-1,3-Dichloropropene	***** NOT FOUND *****		
43 2-Nitropropane	***** NOT FOUND *****		
44 trans-1,3-Dichloropropene	***** NOT FOUND *****		
45 Toluene	***** NOT FOUND *****		
46 Toluene-d8-surrogate	4.86	109.2 ppb	109 % R
47 1,1,2-Trichloroethane	***** NOT FOUND *****		
48 1,3-Dichloropropane	***** NOT FOUND *****		
49 Ethylmethacrylate	***** NOT FOUND *****		
50 Dibromochloromethane	***** NOT FOUND *****		
51 1,2-Dibromomethane	***** NOT FOUND *****		
52 Tetrachloroethene	***** NOT FOUND *****		
53 Chlorobenzene	***** NOT FOUND *****		
54 Chlorobenzene-d5	7.50	100.0 ppb	ISTD3
55 1,1,1,2-Tetrachloroethane	***** NOT FOUND *****		
56 Ethylbenzene	***** NOT FOUND *****		
57 m/p-xylene	***** NOT FOUND *****		
58 Bromoform	***** NOT FOUND *****		
59 Styrene	***** NOT FOUND *****		
60 o-Xylene	***** NOT FOUND *****		
61 cis-1,4-Dichloro-2-butene	***** NOT FOUND *****		
62 trans-1,4-Dichloro-2-butene	***** NOT FOUND *****		
63 Bromofluorobenzene-surrogate	10.42	92.0 ppb	92 % R
64 1,1,2,2-Tetrachloroethane	***** NOT FOUND *****		
65 1,2,3-Trichloropropane	***** NOT FOUND *****		
66 Isopropylbenzene	***** NOT FOUND *****		
67 Bromobenzene	***** NOT FOUND *****		
68 2-Chlorotoluene	***** NOT FOUND *****		
69 n-Propylbenzene	***** NOT FOUND *****		
70 4-Chlorotoluene	***** NOT FOUND *****		
71 1,3,5-Trimethylbenzene	***** NOT FOUND *****		
72 Pentachloroethane	***** NOT FOUND *****		
73 tert-Butylbenzene	***** NOT FOUND *****		
74 1,2,4-Trimethylbenzene	***** NOT FOUND *****		
75 1,3-Dichlorobenzene	***** NOT FOUND *****		
76 1,4-Dichlorobenzene-d4	13.33	100.0 ppb	ISTD4
77 sec-Butylbenzene	***** NOT FOUND *****		
78 1,4-Dichlorobenzene	***** NOT FOUND *****		
79 p-iso-Propyltoluene	***** NOT FOUND *****		
80 1,2-Dichlorobenzene	***** NOT FOUND *****		
81 n-Butylbenzene	***** NOT FOUND *****		
82 1,2-Bromo-3-chloropropane	***** NOT FOUND *****		
83 Nitobenzene	***** NOT FOUND *****		
84 1,2,4-Trichlorobenzene	***** NOT FOUND *****		
85 Naphthalene	***** NOT FOUND *****		
86 1,2,3-Trichlorobenzene	***** NOT FOUND *****		
87 Hexachlorobutadiene	***** NOT FOUND *****		

END OF REPORT

Qualifiers Not Satisfied

Thu May 25 10:41:01 2006

File : C:\HPCHEM\1\DATA\052406\VOA021.D
Operator : dmt
Acquired : 25 May 2006 10:53 am using AcqMethod VOA0506
Instrument : GCD Syste
Sample Name: SB-15 3-8' H2O
Misc Info :
Vial Number: 1



Calibration Last Updated: Thu May 25 08:53:45 2006
Quantitation Report - Summary

Information from Data File:
File : C:\HPCHEM\1\DATA\052406\VOA021.D
Operator : dmt
Acquired : 25 May 2006 10:53 am using AcqMethod VOA0506
Sample Name: SB-15 3-8' H2O
Misc Info :
Vial Number: 1
CurrentMeth: C:\HPCHEM\1\METHODS\VOA0506R.M

Quantitation Settings:
Non-Reference Peak Window: 10.00 Percent
Correlation Window: 0.02 Minutes
Default Multiplier: 1.00
Default Sample Concentration: 0.00
Peak Type Decoding: m -> Manually re-integrated

Number of Compounds in Database: 87

#	Compound	Ret Time	Concentration	Peak Type
1	Pentafluorobenzene	2.94	100.0 ppb	ISTD1#
2	Dichlorofluoromethane	***** NOT FOUND *****		
3	Chloromethane	***** NOT FOUND *****		
4	Vinyl Chloride	***** NOT FOUND *****		
5	Bromomethane	***** NOT FOUND *****		
6	Chloroethane	***** NOT FOUND *****		
7	Trichlorofluoromethane	***** NOT FOUND *****		
8	Acetonitrile	***** NOT FOUND *****		
9	Diethyl ether	***** NOT FOUND *****		
10	1,1-Dichloroethene	***** NOT FOUND *****		
11	Acrylonitrile	***** NOT FOUND *****		
12	1,1,2-Trichlorotrifluoroethane	***** NOT FOUND *****		
13	Methylene Chloride	2.38	3.6 ppb	# c
14	Iodomethane	***** NOT FOUND *****		
15	trans-1,2-Dichloroethene	***** NOT FOUND *****		
16	MTBE	***** NOT FOUND *****		
17	Allyl chloride	***** NOT FOUND *****		
18	1,1-Dichloroethane	***** NOT FOUND *****		
19	Carbon disulfide	***** NOT FOUND *****		
20	Chloroprene	***** NOT FOUND *****		
21	Methacrylonitrile	***** NOT FOUND *****		
22	cis-1,2-Dichloroethene	***** NOT FOUND *****		
23	Methyl acrylate	***** NOT FOUND *****		
24	Chloroform	***** NOT FOUND *****		
25	2,2-Dichloropropane	***** NOT FOUND *****		
26	Dibromofluoromethane-surrogate	2.86	86.6 ppb	87 1.8
27	Isobutyl alcohol	***** NOT FOUND *****		
28	Tetrahydrofuran	***** NOT FOUND *****		
29	1,1,1-Trichloroethane	***** NOT FOUND *****		
30	1,2-Dichloroethane	***** NOT FOUND *****		
31	1,1-Dichloropropene	***** NOT FOUND *****		
32	Benzene	***** NOT FOUND *****		

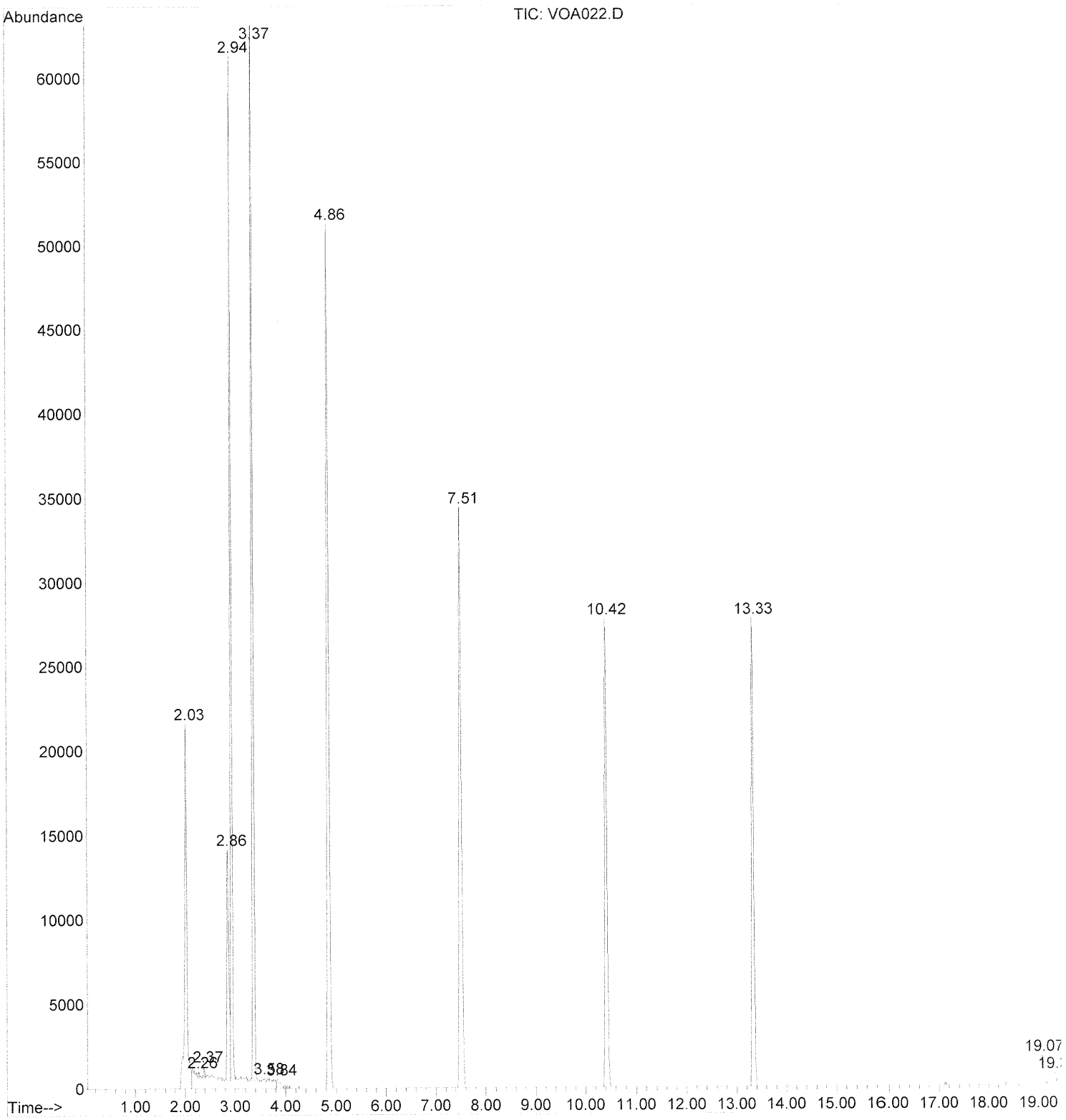
33 Carbon Tetrachloride	***** NOT FOUND *****		
34 2-Chloroethanol	***** NOT FOUND *****		
35 1,4-Difluorobenzene	3.37	100.0 ppb	ISTD2
36 1,2-Dichloropropane	***** NOT FOUND *****		
37 Trichloroethene	***** NOT FOUND *****		
38 Dibromomethane	***** NOT FOUND *****		
39 Bromodichloromethane	***** NOT FOUND *****		
40 Methyl methacrylate	***** NOT FOUND *****		
41 1,4-Dioxane	***** NOT FOUND *****		
42 cis-1,3-Dichloropropene	***** NOT FOUND *****		
43 2-Nitropropane	***** NOT FOUND *****		
44 trans-1,3-Dichloropropene	***** NOT FOUND *****		
45 Toluene	***** NOT FOUND *****		
46 Toluene-d8-surrogate	4.86	99.9 ppb	100 %R
47 1,1,2-Trichloroethane	***** NOT FOUND *****		
48 1,3-Dichloropropane	***** NOT FOUND *****		
49 Ethylmethacrylate	***** NOT FOUND *****		
50 Dibromochloromethane	***** NOT FOUND *****		
51 1,2-Dibromomethane	***** NOT FOUND *****		
52 Tetrachloroethene	***** NOT FOUND *****		
53 Chlorobenzene	***** NOT FOUND *****		
54 Chlorobenzene-d5	7.51	100.0 ppb	ISTD3
55 1,1,1,2-Tetrachloroethane	***** NOT FOUND *****		
56 Ethylbenzene	***** NOT FOUND *****		
57 m/p-xylene	8.45	1.9 ppb	# on D
58 Bromoform	***** NOT FOUND *****		
59 Styrene	***** NOT FOUND *****		
60 o-Xylene	***** NOT FOUND *****		
61 cis-1,4-Dichloro-2-butene	***** NOT FOUND *****		
62 trans-1,4-Dichloro-2-butene	***** NOT FOUND *****		
63 Bromofluorobenzene-surrogate	10.42	96.5 ppb	97 %R
64 1,1,2,2-Tetrachloroethane	***** NOT FOUND *****		
65 1,2,3-Trichloropropane	***** NOT FOUND *****		
66 Isopropylbenzene	***** NOT FOUND *****		
67 Bromobenzene	***** NOT FOUND *****		
68 2-Chlorotoluene	***** NOT FOUND *****		
69 n-Propylbenzene	***** NOT FOUND *****		
70 4-Chlorotoluene	***** NOT FOUND *****		
71 1,3,5-Trimethylbenzene	***** NOT FOUND *****		
72 Pentachloroethane	***** NOT FOUND *****		
73 tert-Butylbenzene	***** NOT FOUND *****		
74 1,2,4-Trimethylbenzene	***** NOT FOUND *****		
75 1,3-Dichlorobenzene	***** NOT FOUND *****		
76 1,4-Dichlorobenzene-d4	13.34	100.0 ppb	ISTD4
77 sec-Butylbenzene	***** NOT FOUND *****		
78 1,4-Dichlorobenzene	***** NOT FOUND *****		
79 p-iso-Propyltoluene	***** NOT FOUND *****		
80 1,2-Dichlorobenzene	***** NOT FOUND *****		
81 n-Butylbenzene	***** NOT FOUND *****		
82 1,2-Bromo-3-chloropropane	***** NOT FOUND *****		
83 Nitobenzene	***** NOT FOUND *****		
84 1,2,4-Trichlorobenzene	***** NOT FOUND *****		
85 Naphthalene	***** NOT FOUND *****		
86 1,2,3-Trichlorobenzene	***** NOT FOUND *****		
87 Hexachlorobutadiene	***** NOT FOUND *****		

END OF REPORT

Qualifiers Not Satisfied

Thu May 25 11:15:36 2006

File : C:\HPCHEM\1\DATA\052406\VOA022.D
Operator : dmt
Acquired : 25 May 2006 11:27 am using AcqMethod VOA0506
Instrument : GCD Syste
Sample Name: SB-26 8-12' H2O
Misc Info :
Vial Number: 1



Calibration Last Updated: Thu May 25 08:53:45 2006
Quantitation Report - Summary

Information from Data File:
File : C:\HPCHEM\1\DATA\052406\VOA022.D
Operator : dmt
Acquired : 25 May 2006 11:27 am using AcqMethod VOA0506
Sample Name: SB-26 8-12' H2O
Misc Info :
Vial Number: 1
CurrentMeth: C:\HPCHEM\1\METHODS\VOA0506R.M

Quantitation Settings:
Non-Reference Peak Window: 10.00 Percent
Correlation Window: 0.02 Minutes
Default Multiplier: 1.00
Default Sample Concentration: 0.00
Peak Type Decoding: m -> Manually re-integrated

Number of Compounds in Database: 87

#	Compound	Ret Time	Concentration	Peak Type
1	Pentafluorobenzene	2.94	100.0 ppb	ISTD1#
2	Dichlorofluoromethane	***** NOT FOUND *****		
3	Chloromethane	***** NOT FOUND *****		
4	Vinyl Chloride	***** NOT FOUND *****		
5	Bromomethane	***** NOT FOUND *****		
6	Chloroethane	***** NOT FOUND *****		
7	Trichlorofluoromethane	***** NOT FOUND *****		
8	Acetonitrile	***** NOT FOUND *****		
9	Diethyl ether	***** NOT FOUND *****		
10	1,1-Dichloroethene	***** NOT FOUND *****		
11	Acrylonitrile	***** NOT FOUND *****		
12	1,1,2-Trichlorotrifluoroethane	***** NOT FOUND *****		
13	Methylene Chloride	2.37	3.0 ppb	#
14	Iodomethane	***** NOT FOUND *****		
15	trans-1,2-Dichloroethene	***** NOT FOUND *****		
16	MTBE	***** NOT FOUND *****		
17	Allyl chloride	***** NOT FOUND *****		
18	1,1-Dichloroethane	***** NOT FOUND *****		
19	Carbon disulfide	***** NOT FOUND *****		
20	Chloroprene	***** NOT FOUND *****		
21	Methacrylonitrile	***** NOT FOUND *****		
22	cis-1,2-Dichloroethene	***** NOT FOUND *****		
23	Methyl acrylate	***** NOT FOUND *****		
24	Chloroform	***** NOT FOUND *****		
25	2,2-Dichloropropane	***** NOT FOUND *****		
26	Dibromofluoromethane-surrogate	2.86	87.9 ppb	
27	Isobutyl alcohol	***** NOT FOUND *****		
28	Tetrahydrofuran	***** NOT FOUND *****		
29	1,1,1-Trichloroethane	***** NOT FOUND *****		
30	1,2-Dichloroethane	***** NOT FOUND *****		
31	1,1-Dichloropropene	***** NOT FOUND *****		
32	Benzene	***** NOT FOUND *****		

88 %R

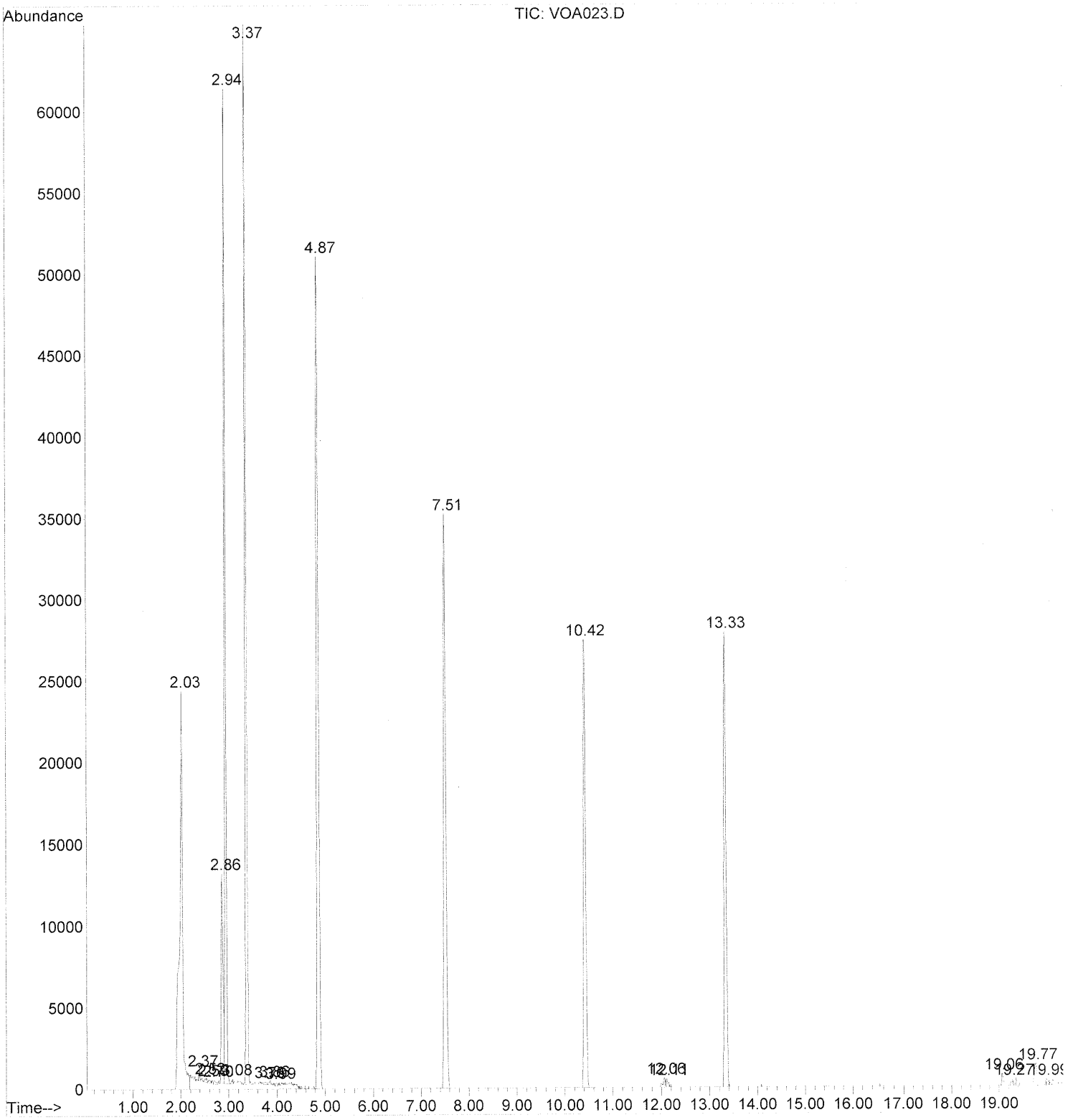
33	Carbon Tetrachloride	***** NOT FOUND *****		
34	2-Chloroethanol	***** NOT FOUND *****		
35	1,4-Difluorobenzene	3.37	100.0 ppb	ISTD2
36	1,2-Dichloropropane	***** NOT FOUND *****		
37	Trichloroethene	***** NOT FOUND *****		
38	Dibromomethane	***** NOT FOUND *****		
39	Bromodichloromethane	***** NOT FOUND *****		
40	Methyl methacrylate	***** NOT FOUND *****		
41	1,4-Dioxane	***** NOT FOUND *****		
42	cis-1,3-Dichloropropene	***** NOT FOUND *****		
43	2-Nitropropane	***** NOT FOUND *****		
44	trans-1,3-Dichloropropene	***** NOT FOUND *****		
45	Toluene	***** NOT FOUND *****		
46	Toluene-d8-surrogate	4.86	99.7 ppb	100 %R
47	1,1,2-Trichloroethane	***** NOT FOUND *****		
48	1,3-Dichloropropane	***** NOT FOUND *****		
49	Ethylmethacrylate	***** NOT FOUND *****		
50	Dibromochloromethane	***** NOT FOUND *****		
51	1,2-Dibromomethane	***** NOT FOUND *****		
52	Tetrachloroethene	***** NOT FOUND *****		
53	Chlorobenzene	***** NOT FOUND *****		
54	Chlorobenzene-d5	7.50	100.0 ppb	ISTD3
55	1,1,1,2-Tetrachloroethane	***** NOT FOUND *****		
56	Ethylbenzene	***** NOT FOUND *****		
57	m/p-xylene	***** NOT FOUND *****		
58	Bromoform	***** NOT FOUND *****		
59	Styrene	***** NOT FOUND *****		
60	o-Xylene	***** NOT FOUND *****		
61	cis-1,4-Dichloro-2-butene	***** NOT FOUND *****		
62	trans-1,4-Dichloro-2-butene	***** NOT FOUND *****		
63	Bromofluorobenzene-surrogate	10.42	98.3 ppb	98 %R
64	1,1,2,2-Tetrachloroethane	***** NOT FOUND *****		
65	1,2,3-Trichloropropane	***** NOT FOUND *****		
66	Isopropylbenzene	***** NOT FOUND *****		
67	Bromobenzene	***** NOT FOUND *****		
68	2-Chlorotoluene	***** NOT FOUND *****		
69	n-Propylbenzene	***** NOT FOUND *****		
70	4-Chlorotoluene	***** NOT FOUND *****		
71	1,3,5-Trimethylbenzene	***** NOT FOUND *****		
72	Pentachloroethane	***** NOT FOUND *****		
73	tert-Butylbenzene	***** NOT FOUND *****		
74	1,2,4-Trimethylbenzene	***** NOT FOUND *****		
75	1,3-Dichlorobenzene	***** NOT FOUND *****		
76	1,4-Dichlorobenzene-d4	13.33	100.0 ppb	ISTD4
77	sec-Butylbenzene	***** NOT FOUND *****		
78	1,4-Dichlorobenzene	***** NOT FOUND *****		
79	p-iso-Propyltoluene	***** NOT FOUND *****		
80	1,2-Dichlorobenzene	***** NOT FOUND *****		
81	n-Butylbenzene	***** NOT FOUND *****		
82	1,2-Bromo-3-chloropropane	***** NOT FOUND *****		
83	Nitobenzene	***** NOT FOUND *****		
84	1,2,4-Trichlorobenzene	***** NOT FOUND *****		
85	Naphthalene	***** NOT FOUND *****		
86	1,2,3-Trichlorobenzene	***** NOT FOUND *****		
87	Hexachlorobutadiene	***** NOT FOUND *****		

END OF REPORT

Qualifiers Not Satisfied

Thu May 25 11:47:37 2006

File : C:\HPCHEM\1\DATA\052406\VOA023.D
Operator : dmt
Acquired : 25 May 2006 11:53 am using AcqMethod VOA0506
Instrument : GCD Syste
Sample Name: SB-25 7-11' H2O
Misc Info :
Vial Number: 1



Calibration Last Updated: Thu May 25 08:53:45 2006
Quantitation Report - Summary

Information from Data File:
File : C:\HPCHEM\1\DATA\052406\VOA023.D
Operator : dmt
Acquired : 25 May 2006 11:53 am using AcqMethod VOA0506
Sample Name: SB-25 7-11' H2O
Misc Info :
Vial Number: 1
CurrentMeth: C:\HPCHEM\1\METHODS\VOA0506R.M

Quantitation Settings:
Non-Reference Peak Window: 10.00 Percent
Correlation Window: 0.02 Minutes
Default Multiplier: 1.00
Default Sample Concentration: 0.00
Peak Type Decoding: m -> Manually re-integrated

Number of Compounds in Database: 87

#	Compound	Ret Time	Concentration	Peak Type
1	Pentafluorobenzene	2.94	100.0 ppb	ISTD1
2	Dichlorofluoromethane	***** NOT FOUND *****		
3	Chloromethane	***** NOT FOUND *****		
4	Vinyl Chloride	***** NOT FOUND *****		
5	Bromomethane	***** NOT FOUND *****		
6	Chloroethane	***** NOT FOUND *****		
7	Trichlorofluoromethane	***** NOT FOUND *****		
8	Acetonitrile	***** NOT FOUND *****		
9	Diethyl ether	***** NOT FOUND *****		
10	1,1-Dichloroethene	***** NOT FOUND *****		
11	Acrylonitrile	***** NOT FOUND *****		
12	1,1,2-Trichlorotrifluoroethane	***** NOT FOUND *****		
13	Methylene Chloride	2.37	3.3 ppb	#
14	Iodomethane	***** NOT FOUND *****		
15	trans-1,2-Dichloroethene	***** NOT FOUND *****		
16	MTBE	***** NOT FOUND *****		
17	Allyl chloride	***** NOT FOUND *****		
18	1,1-Dichloroethane	***** NOT FOUND *****		
19	Carbon disulfide	***** NOT FOUND *****		
20	Chloroprene	***** NOT FOUND *****		
21	Methacrylonitrile	***** NOT FOUND *****		
22	cis-1,2-Dichloroethene	***** NOT FOUND *****		
23	Methyl acrylate	***** NOT FOUND *****		
24	Chloroform	***** NOT FOUND *****		
25	2,2-Dichloropropane	***** NOT FOUND *****		
26	Dibromofluoromethane-surrogate	2.86	87.0 ppb	
27	Isobutyl alcohol	***** NOT FOUND *****		
28	Tetrahydrofuran	***** NOT FOUND *****		
29	1,1,1-Trichloroethane	***** NOT FOUND *****		
30	1,2-Dichloroethane	***** NOT FOUND *****		
31	1,1-Dichloropropene	***** NOT FOUND *****		
32	Benzene	***** NOT FOUND *****		

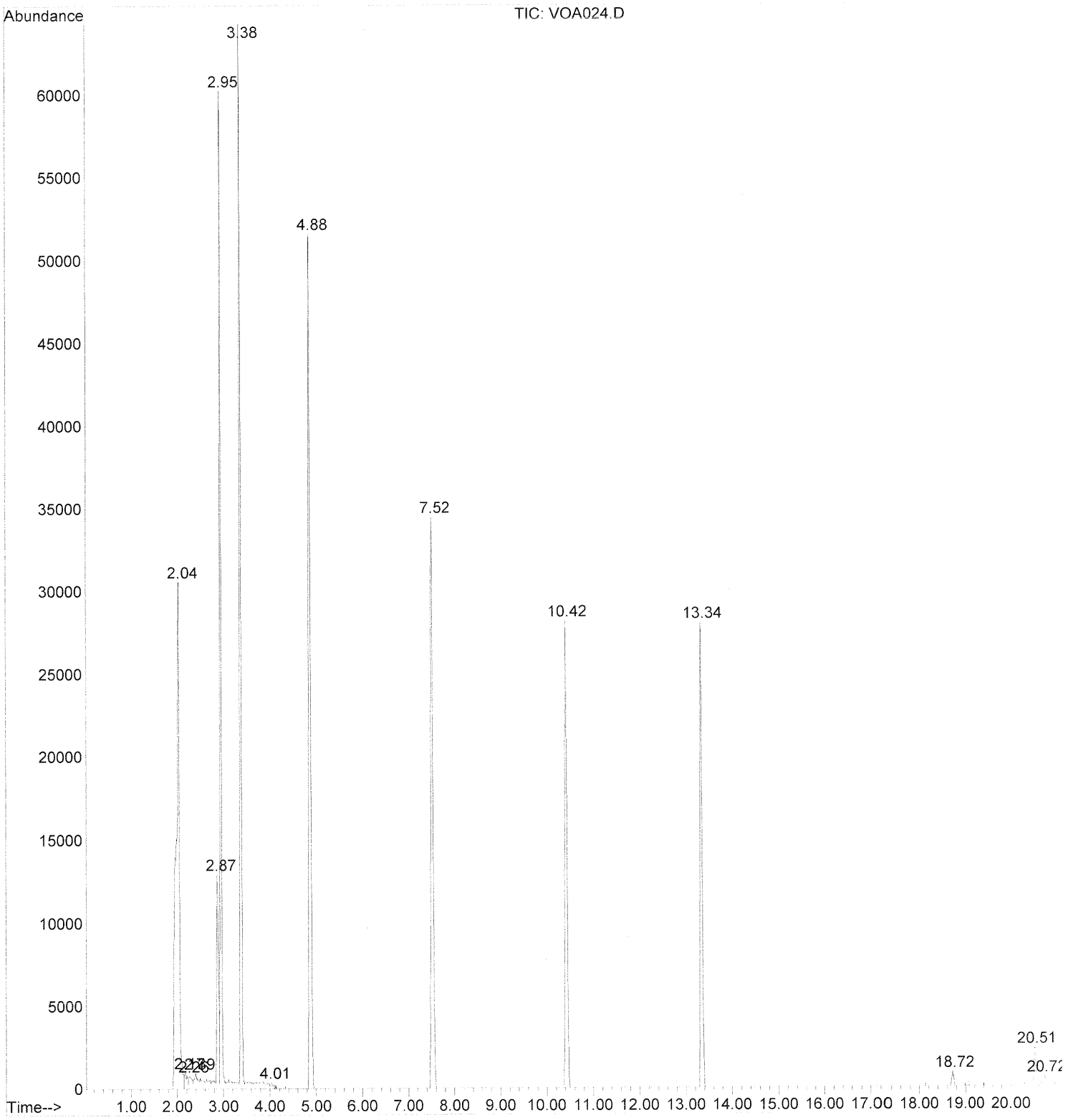
33	Carbon Tetrachloride	***** NOT FOUND *****		
34	2-Chloroethanol	***** NOT FOUND *****		
35	1,4-Difluorobenzene	3.37	100.0 ppb	ISTD2
36	1,2-Dichloropropane	***** NOT FOUND *****		
37	Trichloroethene	***** NOT FOUND *****		
38	Dibromomethane	***** NOT FOUND *****		
39	Bromodichloromethane	***** NOT FOUND *****		
40	Methyl methacrylate	***** NOT FOUND *****		
41	1,4-Dioxane	***** NOT FOUND *****		
42	cis-1,3-Dichloropropene	***** NOT FOUND *****		
43	2-Nitropropane	***** NOT FOUND *****		
44	trans-1,3-Dichloropropene	***** NOT FOUND *****		
45	Toluene	***** NOT FOUND *****		
46	Toluene-d8-surrogate	4.87	106.5 ppb	107 %R
47	1,1,2-Trichloroethane	***** NOT FOUND *****		
48	1,3-Dichloropropane	***** NOT FOUND *****		
49	Ethylmethacrylate	***** NOT FOUND *****		
50	Dibromochloromethane	***** NOT FOUND *****		
51	1,2-Dibromomethane	***** NOT FOUND *****		
52	Tetrachloroethene	***** NOT FOUND *****		
53	Chlorobenzene	***** NOT FOUND *****		
54	Chlorobenzene-d5	7.51	100.0 ppb	ISTD3
55	1,1,1,2-Tetrachloroethane	***** NOT FOUND *****		
56	Ethylbenzene	***** NOT FOUND *****		
57	m/p-xylene	***** NOT FOUND *****		
58	Bromoform	***** NOT FOUND *****		
59	Styrene	***** NOT FOUND *****		
60	o-Xylene	***** NOT FOUND *****		
61	cis-1,4-Dichloro-2-butene	***** NOT FOUND *****		
62	trans-1,4-Dichloro-2-butene	***** NOT FOUND *****		
63	Bromofluorobenzene-surrogate	10.43	92.2 ppb	92 %R
64	1,1,2,2-Tetrachloroethane	***** NOT FOUND *****		
65	1,2,3-Trichloropropane	***** NOT FOUND *****		
66	Isopropylbenzene	***** NOT FOUND *****		
67	Bromobenzene	***** NOT FOUND *****		
68	2-Chlorotoluene	***** NOT FOUND *****		
69	n-Propylbenzene	***** NOT FOUND *****		
70	4-Chlorotoluene	***** NOT FOUND *****		
71	1,3,5-Trimethylbenzene	***** NOT FOUND *****		
72	Pentachloroethane	***** NOT FOUND *****		
73	tert-Butylbenzene	***** NOT FOUND *****		
74	1,2,4-Trimethylbenzene	***** NOT FOUND *****		
75	1,3-Dichlorobenzene	***** NOT FOUND *****		
76	1,4-Dichlorobenzene-d4	13.33	100.0 ppb	ISTD4
77	sec-Butylbenzene	***** NOT FOUND *****		
78	1,4-Dichlorobenzene	***** NOT FOUND *****		
79	p-iso-Propyltoluene	***** NOT FOUND *****		
80	1,2-Dichlorobenzene	***** NOT FOUND *****		
81	n-Butylbenzene	***** NOT FOUND *****		
82	1,2-Bromo-3-chloropropane	***** NOT FOUND *****		
83	Nitobenzene	***** NOT FOUND *****		
84	1,2,4-Trichlorobenzene	***** NOT FOUND *****		
85	Naphthalene	***** NOT FOUND *****		
86	1,2,3-Trichlorobenzene	***** NOT FOUND *****		
87	Hexachlorobutadiene	***** NOT FOUND *****		

END OF REPORT

Qualifiers Not Satisfied

Thu May 25 12:14:38 2006

File : C:\HPCHEM\1\DATA\052406\VOA024.D
Operator : dmt
Acquired : 25 May 2006 12:20 pm using AcqMethod VOA0506
Instrument : GCD Syste
Sample Name: SB-30 6-10' H2O
Misc Info :
Vial Number: 1



Calibration Last Updated: Thu May 25 08:53:45 2006
Quantitation Report - Summary

Information from Data File:
File : C:\HPCHEM\1\DATA\052406\VOA024.D
Operator : dmt
Acquired : 25 May 2006 12:20 pm using AcqMethod VOA0506
Sample Name: SB-30 6-10' H2O
Misc Info :
Vial Number: 1
CurrentMeth: C:\HPCHEM\1\METHODS\VOA0506R.M

Quantitation Settings:
Non-Reference Peak Window: 10.00 Percent
Correlation Window: 0.02 Minutes
Default Multiplier: 1.00
Default Sample Concentration: 0.00
Peak Type Decoding: m -> Manually re-integrated

Number of Compounds in Database: 87

#	Compound	Ret Time	Concentration	Peak Type
1	Pentafluorobenzene	2.95	100.0 ppb	ISTD1
2	Dichlorofluoromethane	***** NOT FOUND *****		
3	Chloromethane	***** NOT FOUND *****		
4	Vinyl Chloride	***** NOT FOUND *****		
5	Bromomethane	***** NOT FOUND *****		
6	Chloroethane	***** NOT FOUND *****		
7	Trichlorofluoromethane	***** NOT FOUND *****		
8	Acetonitrile	***** NOT FOUND *****		
9	Diethyl ether	***** NOT FOUND *****		
10	1,1-Dichloroethene	***** NOT FOUND *****		
11	Acrylonitrile	***** NOT FOUND *****		
12	1,1,2-Trichlorotrifluoroethane	***** NOT FOUND *****		
13	Methylene Chloride	2.39	4.9 ppb	#
14	Iodomethane	***** NOT FOUND *****		
15	trans-1,2-Dichloroethene	***** NOT FOUND *****		
16	MTBE	***** NOT FOUND *****		
17	Allyl chloride	***** NOT FOUND *****		
18	1,1-Dichloroethane	***** NOT FOUND *****		
19	Carbon disulfide	***** NOT FOUND *****		
20	Chloroprene	***** NOT FOUND *****		
21	Methacrylonitrile	***** NOT FOUND *****		
22	cis-1,2-Dichloroethene	***** NOT FOUND *****		
23	Methyl acrylate	***** NOT FOUND *****		
24	Chloroform	***** NOT FOUND *****		
25	2,2-Dichloropropane	***** NOT FOUND *****		
26	Dibromofluoromethane-surrogate	2.87	90.3 ppb	
27	Isobutyl alcohol	***** NOT FOUND *****		
28	Tetrahydrofuran	***** NOT FOUND *****		
29	1,1,1-Trichloroethane	***** NOT FOUND *****		
30	1,2-Dichloroethane	***** NOT FOUND *****		
31	1,1-Dichloropropene	***** NOT FOUND *****		
32	Benzene	***** NOT FOUND *****		

90 %R

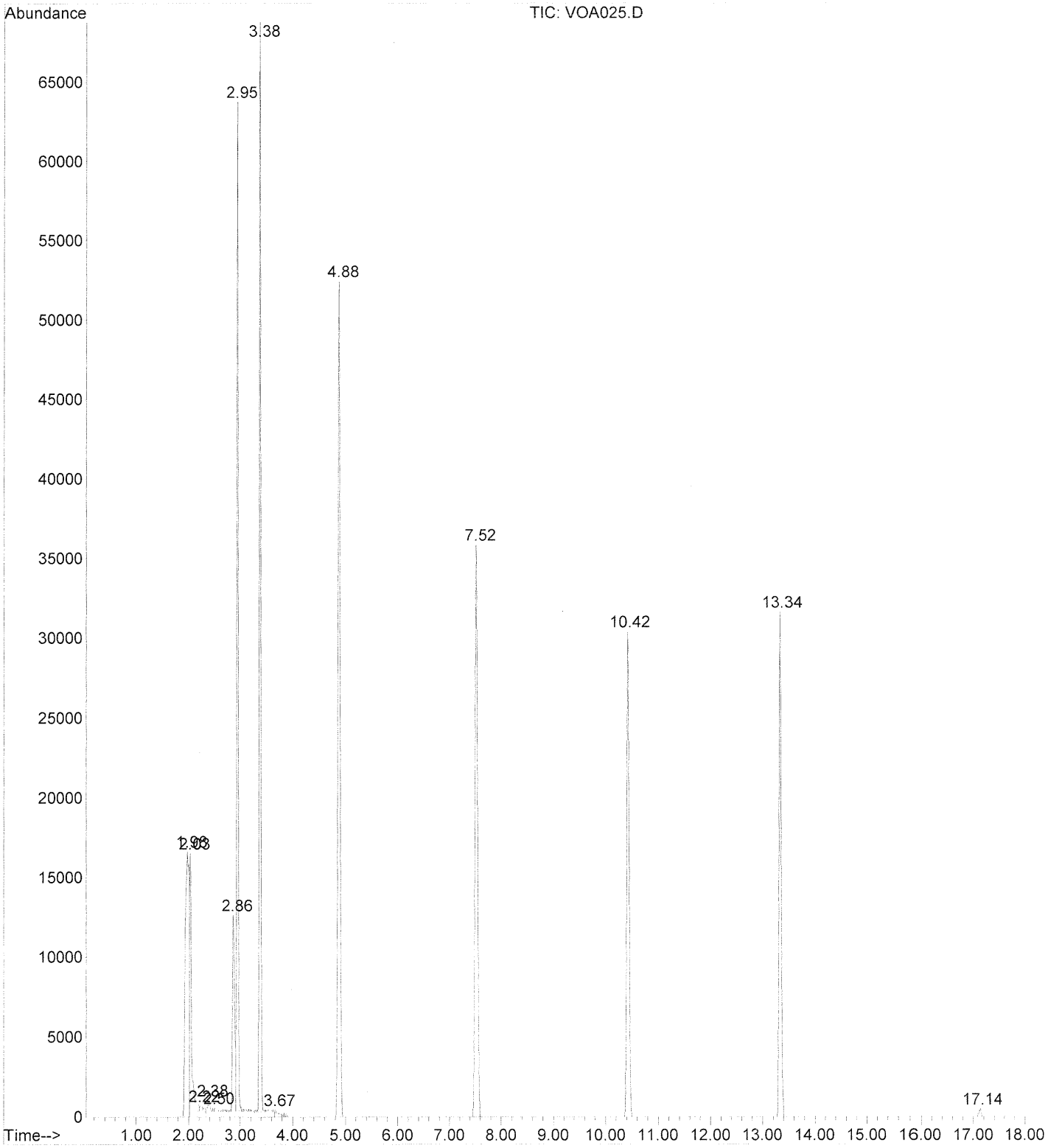
33	Carbon Tetrachloride	***** NOT FOUND *****		
34	2-Chloroethanol	***** NOT FOUND *****		
35	1,4-Difluorobenzene	3.38	100.0 ppb	ISTD2
36	1,2-Dichloropropane	***** NOT FOUND *****		
37	Trichloroethene	***** NOT FOUND *****		
38	Dibromomethane	***** NOT FOUND *****		
39	Bromodichloromethane	***** NOT FOUND *****		
40	Methyl methacrylate	***** NOT FOUND *****		
41	1,4-Dioxane	***** NOT FOUND *****		
42	cis-1,3-Dichloropropene	***** NOT FOUND *****		
43	2-Nitropropane	***** NOT FOUND *****		
44	trans-1,3-Dichloropropene	***** NOT FOUND *****		
45	Toluene	***** NOT FOUND *****		
46	Toluene-d8-surrogate	4.88	107.1 ppb	107% R
47	1,1,2-Trichloroethane	***** NOT FOUND *****		
48	1,3-Dichloropropane	***** NOT FOUND *****		
49	Ethylmethacrylate	***** NOT FOUND *****		
50	Dibromochloromethane	***** NOT FOUND *****		
51	1,2-Dibromomethane	***** NOT FOUND *****		
52	Tetrachloroethene	***** NOT FOUND *****		
53	Chlorobenzene	***** NOT FOUND *****		
54	Chlorobenzene-d5	7.53	100.0 ppb	ISTD3
55	1,1,1,2-Tetrachloroethane	***** NOT FOUND *****		
56	Ethylbenzene	***** NOT FOUND *****		
57	m/p-xylene	***** NOT FOUND *****		
58	Bromoform	***** NOT FOUND *****		
59	Styrene	***** NOT FOUND *****		
60	o-Xylene	***** NOT FOUND *****		
61	cis-1,4-Dichloro-2-butene	***** NOT FOUND *****		
62	trans-1,4-Dichloro-2-butene	***** NOT FOUND *****		
63	Bromofluorobenzene-surrogate	10.42	92.6 ppb	93% R
64	1,1,2,2-Tetrachloroethane	***** NOT FOUND *****		
65	1,2,3-Trichloropropane	***** NOT FOUND *****		
66	Isopropylbenzene	***** NOT FOUND *****		
67	Bromobenzene	***** NOT FOUND *****		
68	2-Chlorotoluene	***** NOT FOUND *****		
69	n-Propylbenzene	***** NOT FOUND *****		
70	4-Chlorotoluene	***** NOT FOUND *****		
71	1,3,5-Trimethylbenzene	***** NOT FOUND *****		
72	Pentachloroethane	***** NOT FOUND *****		
73	tert-Butylbenzene	***** NOT FOUND *****		
74	1,2,4-Trimethylbenzene	***** NOT FOUND *****		
75	1,3-Dichlorobenzene	***** NOT FOUND *****		
76	1,4-Dichlorobenzene-d4	13.34	100.0 ppb	ISTD4
77	sec-Butylbenzene	***** NOT FOUND *****		
78	1,4-Dichlorobenzene	***** NOT FOUND *****		
79	p-iso-Propyltoluene	***** NOT FOUND *****		
80	1,2-Dichlorobenzene	***** NOT FOUND *****		
81	n-Butylbenzene	***** NOT FOUND *****		
82	1,2-Bromo-3-chloropropane	***** NOT FOUND *****		
83	Nitobenzene	***** NOT FOUND *****		
84	1,2,4-Trichlorobenzene	***** NOT FOUND *****		
85	Naphthalene	***** NOT FOUND *****		
86	1,2,3-Trichlorobenzene	***** NOT FOUND *****		
87	Hexachlorobutadiene	***** NOT FOUND *****		

END OF REPORT

Qualifiers Not Satisfied

Thu May 25 12:46:56 2006

File : C:\HPCHEM\1\DATA\052406\VOA025.D
Operator : dmt
Acquired : 25 May 2006 12:54 pm using AcqMethod VOA0506
Instrument : GCD Syste
Sample Name: SB-31 4-8' H2O
Misc Info :
Vial Number: 1



Calibration Last Updated: Thu May 25 08:53:45 2006
Quantitation Report - Summary

Information from Data File:
File : C:\HPCHEM\1\DATA\052406\VOA025.D
Operator : dmt
Acquired : 25 May 2006 12:54 pm using AcqMethod VOA0506
Sample Name: SB-31 4-8' H2O
Misc Info :
Vial Number: 1
CurrentMeth: C:\HPCHEM\1\METHODS\VOA0506R.M

Quantitation Settings:
Non-Reference Peak Window: 10.00 Percent
Correlation Window: 0.02 Minutes
Default Multiplier: 1.00
Default Sample Concentration: 0.00
Peak Type Decoding: m -> Manually re-integrated

Number of Compounds in Database: 87

#	Compound	Ret Time	Concentration	Peak Type
1	Pentafluorobenzene	2.95	100.0 ppb	ISTD1
2	Dichlorofluoromethane	***** NOT FOUND *****		
3	Chloromethane	***** NOT FOUND *****		
4	Vinyl Chloride	***** NOT FOUND *****		
5	Bromomethane	***** NOT FOUND *****		
6	Chloroethane	***** NOT FOUND *****		
7	Trichlorofluoromethane	***** NOT FOUND *****		
8	Acetonitrile	***** NOT FOUND *****		
9	Diethyl ether	***** NOT FOUND *****		
10	1,1-Dichloroethene	***** NOT FOUND *****		
11	Acrylonitrile	***** NOT FOUND *****		
12	1,1,2-Trichlorotrifluoroethane	***** NOT FOUND *****		
13	Methylene Chloride	2.38	4.8 ppb	#
14	Iodomethane	***** NOT FOUND *****		
15	trans-1,2-Dichloroethene	***** NOT FOUND *****		
16	MTBE	***** NOT FOUND *****		
17	Allyl chloride	***** NOT FOUND *****		
18	1,1-Dichloroethane	***** NOT FOUND *****		
19	Carbon disulfide	***** NOT FOUND *****		
20	Chloroprene	***** NOT FOUND *****		
21	Methacrylonitrile	***** NOT FOUND *****		
22	cis-1,2-Dichloroethene	***** NOT FOUND *****		
23	Methyl acrylate	***** NOT FOUND *****		
24	Chloroform	***** NOT FOUND *****		
25	2,2-Dichloropropane	***** NOT FOUND *****		
26	Dibromofluoromethane-surrogate	2.87	91.3 ppb	
27	Isobutyl alcohol	***** NOT FOUND *****		
28	Tetrahydrofuran	***** NOT FOUND *****		
29	1,1,1-Trichloroethane	***** NOT FOUND *****		
30	1,2-Dichloroethane	***** NOT FOUND *****		
31	1,1-Dichloropropene	***** NOT FOUND *****		
32	Benzene	***** NOT FOUND *****		

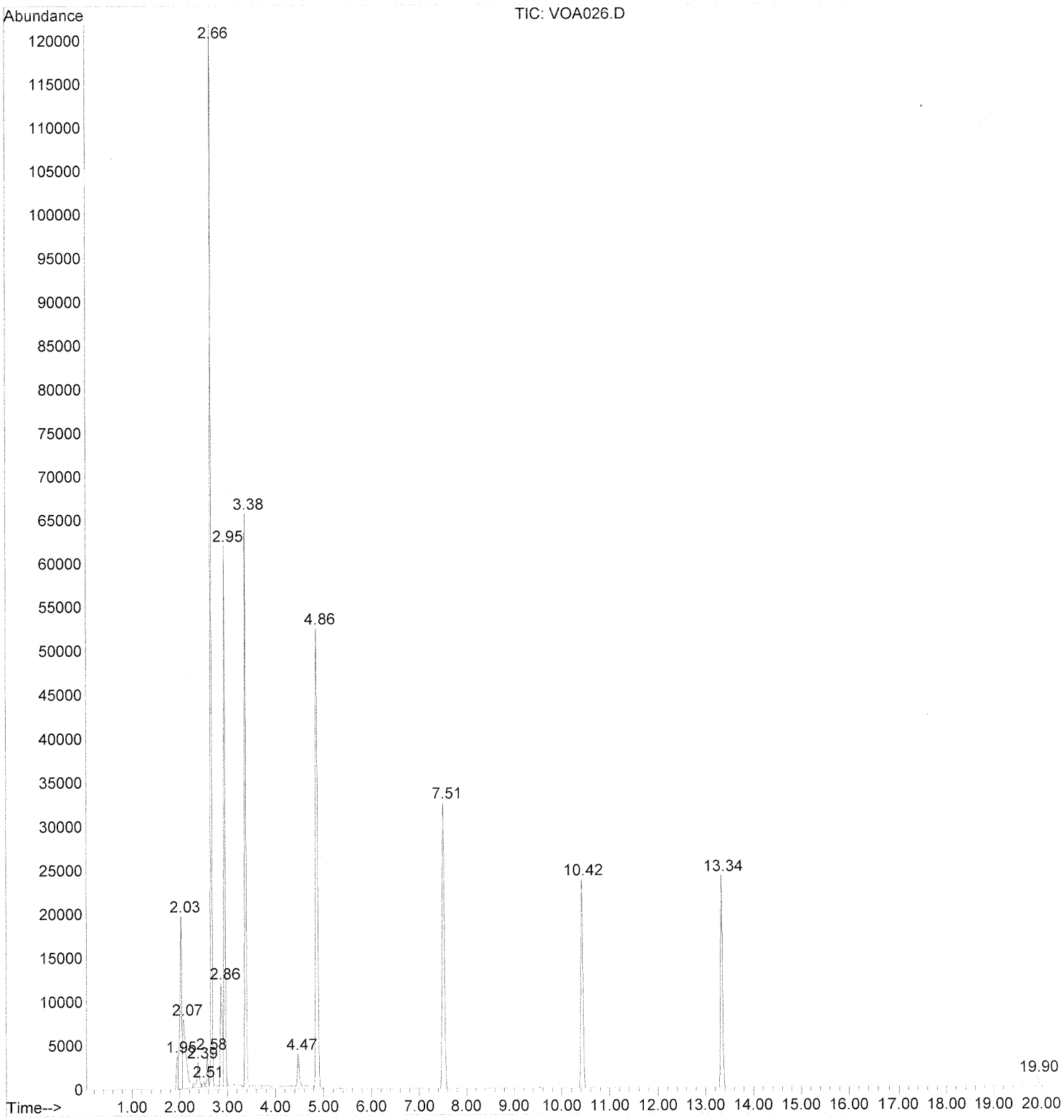
33 Carbon Tetrachloride	***** NOT FOUND *****		
34 2-Chloroethanol	***** NOT FOUND *****		
35 1,4-Difluorobenzene	3.38	100.0 ppb	ISTD2
36 1,2-Dichloropropane	***** NOT FOUND *****		
37 Trichloroethene	***** NOT FOUND *****		
38 Dibromomethane	***** NOT FOUND *****		
39 Bromodichloromethane	***** NOT FOUND *****		
40 Methyl methacrylate	***** NOT FOUND *****		
41 1,4-Dioxane	***** NOT FOUND *****		
42 cis-1,3-Dichloropropene	***** NOT FOUND *****		
43 2-Nitropropane	***** NOT FOUND *****		
44 trans-1,3-Dichloropropene	***** NOT FOUND *****		
45 Toluene	***** NOT FOUND *****		
46 Toluene-d8-surrogate	4.88	100.5 ppb	
47 1,1,2-Trichloroethane	***** NOT FOUND *****		
48 1,3-Dichloropropane	***** NOT FOUND *****		
49 Ethylmethacrylate	***** NOT FOUND *****		
50 Dibromochloromethane	***** NOT FOUND *****		
51 1,2-Dibromomethane	***** NOT FOUND *****		
52 Tetrachloroethene	***** NOT FOUND *****		
53 Chlorobenzene	***** NOT FOUND *****		
54 Chlorobenzene-d5	7.52	100.0 ppb	ISTD3
55 1,1,1,2-Tetrachloroethane	***** NOT FOUND *****		
56 Ethylbenzene	***** NOT FOUND *****		
57 m/p-xylene	***** NOT FOUND *****		
58 Bromoform	***** NOT FOUND *****		
59 Styrene	***** NOT FOUND *****		
60 o-Xylene	***** NOT FOUND *****		
61 cis-1,4-Dichloro-2-butene	***** NOT FOUND *****		
62 trans-1,4-Dichloro-2-butene	***** NOT FOUND *****		
63 Bromofluorobenzene-surrogate	10.43	96.2 ppb	
64 1,1,2,2-Tetrachloroethane	***** NOT FOUND *****		
65 1,2,3-Trichloropropane	***** NOT FOUND *****		
66 Isopropylbenzene	***** NOT FOUND *****		
67 Bromobenzene	***** NOT FOUND *****		
68 2-Chlorotoluene	***** NOT FOUND *****		
69 n-Propylbenzene	***** NOT FOUND *****		
70 4-Chlorotoluene	***** NOT FOUND *****		
71 1,3,5-Trimethylbenzene	***** NOT FOUND *****		
72 Pentachloroethane	***** NOT FOUND *****		
73 tert-Butylbenzene	***** NOT FOUND *****		
74 1,2,4-Trimethylbenzene	***** NOT FOUND *****		
75 1,3-Dichlorobenzene	***** NOT FOUND *****		
76 1,4-Dichlorobenzene-d4	13.34	100.0 ppb	ISTD4
77 sec-Butylbenzene	***** NOT FOUND *****		
78 1,4-Dichlorobenzene	***** NOT FOUND *****		
79 p-iso-Propyltoluene	***** NOT FOUND *****		
80 1,2-Dichlorobenzene	***** NOT FOUND *****		
81 n-Butylbenzene	***** NOT FOUND *****		
82 1,2-Bromo-3-chloropropane	***** NOT FOUND *****		
83 Nitobenzene	***** NOT FOUND *****		
84 1,2,4-Trichlorobenzene	***** NOT FOUND *****		
85 Naphthalene	***** NOT FOUND *****		
86 1,2,3-Trichlorobenzene	***** NOT FOUND *****		
87 Hexachlorobutadiene	***** NOT FOUND *****		

END OF REPORT

Qualifiers Not Satisfied

Thu May 25 13:14:35 2006

File : C:\HPCHEM\1\DATA\052406\VOA026.D
Operator : dmt
Acquired : 25 May 2006 1:22 pm using AcqMethod VOA0506
Instrument : GCD Syste
Sample Name: SB-27 0-4' SOIL
Misc Info :
Vial Number: 1



Calibration Last Updated: Thu May 25 13:44:00 2006
Quantitation Report - Summary

Information from Data File:
File : C:\HPCHEM\1\DATA\052406\VOA026.D
Operator : dmt
Acquired : 25 May 2006 1:22 pm using AcqMethod VOA0506
Sample Name: SB-27 0-4' SOIL
Misc Info :
Vial Number: 1
CurrentMeth: C:\HPCHEM\1\METHODS\VOA0506R.M

Quantitation Settings:
Non-Reference Peak Window: 10.00 Percent
Correlation Window: 0.02 Minutes
Default Multiplier: 1.00
Default Sample Concentration: 0.00
Peak Type Decoding: m -> Manually re-integrated

Number of Compounds in Database: 87

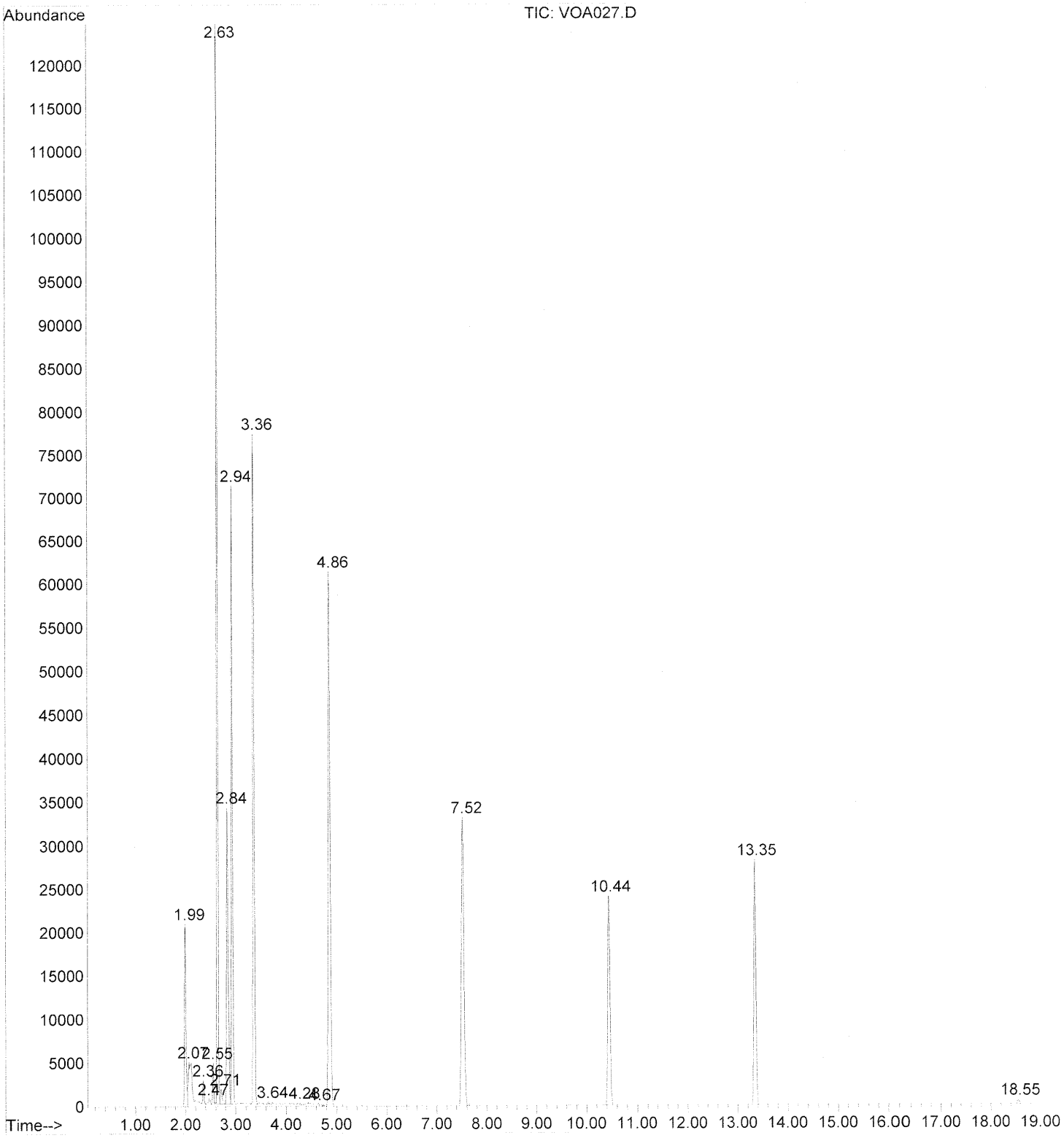
#	Compound	Ret Time	Concentration	Peak Type
1	Pentafluorobenzene	2.95	100.0 ppb	ISTD1
2	Dichlorofluoromethane	***** NOT FOUND *****		
3	Chloromethane	***** NOT FOUND *****		
4	Vinyl Chloride	***** NOT FOUND *****		
5	Bromomethane	***** NOT FOUND *****		
6	Chloroethane	***** NOT FOUND *****		
7	Trichlorofluoromethane	***** NOT FOUND *****		
8	Acetonitrile	***** NOT FOUND *****		
9	Diethyl ether	***** NOT FOUND *****		
10	1,1-Dichloroethene	***** NOT FOUND *****		
11	Acrylonitrile	***** NOT FOUND *****		
12	1,1,2-Trichlorotrifluoroethane	***** NOT FOUND *****		
13	Methylene Chloride	2.39	15.3 ppb	<
14	Iodomethane	***** NOT FOUND *****		
15	trans-1,2-Dichloroethene	***** NOT FOUND *****		
16	MTBE	***** NOT FOUND *****		
17	Allyl chloride	***** NOT FOUND *****		
18	1,1-Dichloroethane	***** NOT FOUND *****		
19	Carbon disulfide	***** NOT FOUND *****		
20	Chloroprene	***** NOT FOUND *****		
21	Methacrylonitrile	***** NOT FOUND *****		
22	cis-1,2-Dichloroethene	***** NOT FOUND *****		
23	Methyl acrylate	***** NOT FOUND *****		
24	Chloroform	***** NOT FOUND *****		
25	2,2-Dichloropropane	***** NOT FOUND *****		
26	Dibromofluoromethane-surrogate	2.86	77.3 ppb	77.1
27	Isobutyl alcohol	***** NOT FOUND *****		
28	Tetrahydrofuran	***** NOT FOUND *****		
29	1,1,1-Trichloroethane	***** NOT FOUND *****		
30	1,2-Dichloroethane	***** NOT FOUND *****		
31	1,1-Dichloropropene	***** NOT FOUND *****		
32	Benzene	***** NOT FOUND *****		

33	Carbon Tetrachloride	***** NOT FOUND *****		
34	2-Chloroethanol	***** NOT FOUND *****		
35	1,4-Difluorobenzene	3.38	100.0 ppb	ISTD2
36	1,2-Dichloropropane	***** NOT FOUND *****		
37	Trichloroethene	***** NOT FOUND *****		
38	Dibromomethane	***** NOT FOUND *****		
39	Bromodichloromethane	***** NOT FOUND *****		
40	Methyl methacrylate	***** NOT FOUND *****		
41	1,4-Dioxane	***** NOT FOUND *****		
42	cis-1,3-Dichloropropene	***** NOT FOUND *****		
43	2-Nitropropane	***** NOT FOUND *****		
44	trans-1,3-Dichloropropene	***** NOT FOUND *****		
45	Toluene	***** NOT FOUND *****		
46	Toluene-d8-surrogate	4.86	111.2 ppb	111 %
47	1,1,2-Trichloroethane	***** NOT FOUND *****		
48	1,3-Dichloropropane	***** NOT FOUND *****		
49	Ethylmethacrylate	***** NOT FOUND *****		
50	Dibromochloromethane	***** NOT FOUND *****		
51	1,2-Dibromomethane	***** NOT FOUND *****		
52	Tetrachloroethene	***** NOT FOUND *****		
53	Chlorobenzene	***** NOT FOUND *****		
54	Chlorobenzene-d5	7.51	100.0 ppb	ISTD3
55	1,1,1,2-Tetrachloroethane	***** NOT FOUND *****		
56	Ethylbenzene	***** NOT FOUND *****		
57	m/p-xylene	***** NOT FOUND *****		
58	Bromoform	***** NOT FOUND *****		
59	Styrene	***** NOT FOUND *****		
60	o-Xylene	***** NOT FOUND *****		
61	cis-1,4-Dichloro-2-butene	***** NOT FOUND *****		
62	trans-1,4-Dichloro-2-butene	***** NOT FOUND *****		
63	Bromofluorobenzene-surrogate	10.42	84.7 ppb	85 %
64	1,1,2,2-Tetrachloroethane	***** NOT FOUND *****		
65	1,2,3-Trichloropropane	***** NOT FOUND *****		
66	Isopropylbenzene	***** NOT FOUND *****		
67	Bromobenzene	***** NOT FOUND *****		
68	2-Chlorotoluene	***** NOT FOUND *****		
69	n-Propylbenzene	***** NOT FOUND *****		
70	4-Chlorotoluene	***** NOT FOUND *****		
71	1,3,5-Trimethylbenzene	***** NOT FOUND *****		
72	Pentachloroethane	***** NOT FOUND *****		
73	tert-Butylbenzene	***** NOT FOUND *****		
74	1,2,4-Trimethylbenzene	***** NOT FOUND *****		
75	1,3-Dichlorobenzene	***** NOT FOUND *****		
76	1,4-Dichlorobenzene-d4	13.34	100.0 ppb	ISTD4
77	sec-Butylbenzene	***** NOT FOUND *****		
78	1,4-Dichlorobenzene	***** NOT FOUND *****		
79	p-iso-Propyltoluene	***** NOT FOUND *****		
80	1,2-Dichlorobenzene	***** NOT FOUND *****		
81	n-Butylbenzene	***** NOT FOUND *****		
82	1,2-Bromo-3-chloropropane	***** NOT FOUND *****		
83	Nitobenzene	***** NOT FOUND *****		
84	1,2,4-Trichlorobenzene	***** NOT FOUND *****		
85	Naphthalene	***** NOT FOUND *****		
86	1,2,3-Trichlorobenzene	***** NOT FOUND *****		
87	Hexachlorobutadiene	***** NOT FOUND *****		

END OF REPORT

Thu May 25 13:44:05 2006

File : C:\HPCHEM\1\DATA\052406\VOA027.D
Operator : dmt
Acquired : 25 May 2006 1:49 pm using AcqMethod VOA0506
Instrument : GCD Syste
Sample Name: SB-29 0-4' SOIL
Misc Info :
Vial Number: 1



Calibration Last Updated: Thu May 25 13:44:00 2006
Quantitation Report - Summary

Information from Data File:
File : C:\HPCHEM\1\DATA\052406\VOA027.D
Operator : dmt
Acquired : 25 May 2006 1:49 pm using AcqMethod VOA0506
Sample Name: SB-29 0-4' SOIL
Misc Info :
Vial Number: 1
CurrentMeth: C:\HPCHEM\1\METHODS\VOA0506R.M

Quantitation Settings:
Non-Reference Peak Window: 10.00 Percent
Correlation Window: 0.02 Minutes
Default Multiplier: 1.00
Default Sample Concentration: 0.00
Peak Type Decoding: m -> Manually re-integrated

Number of Compounds in Database: 87

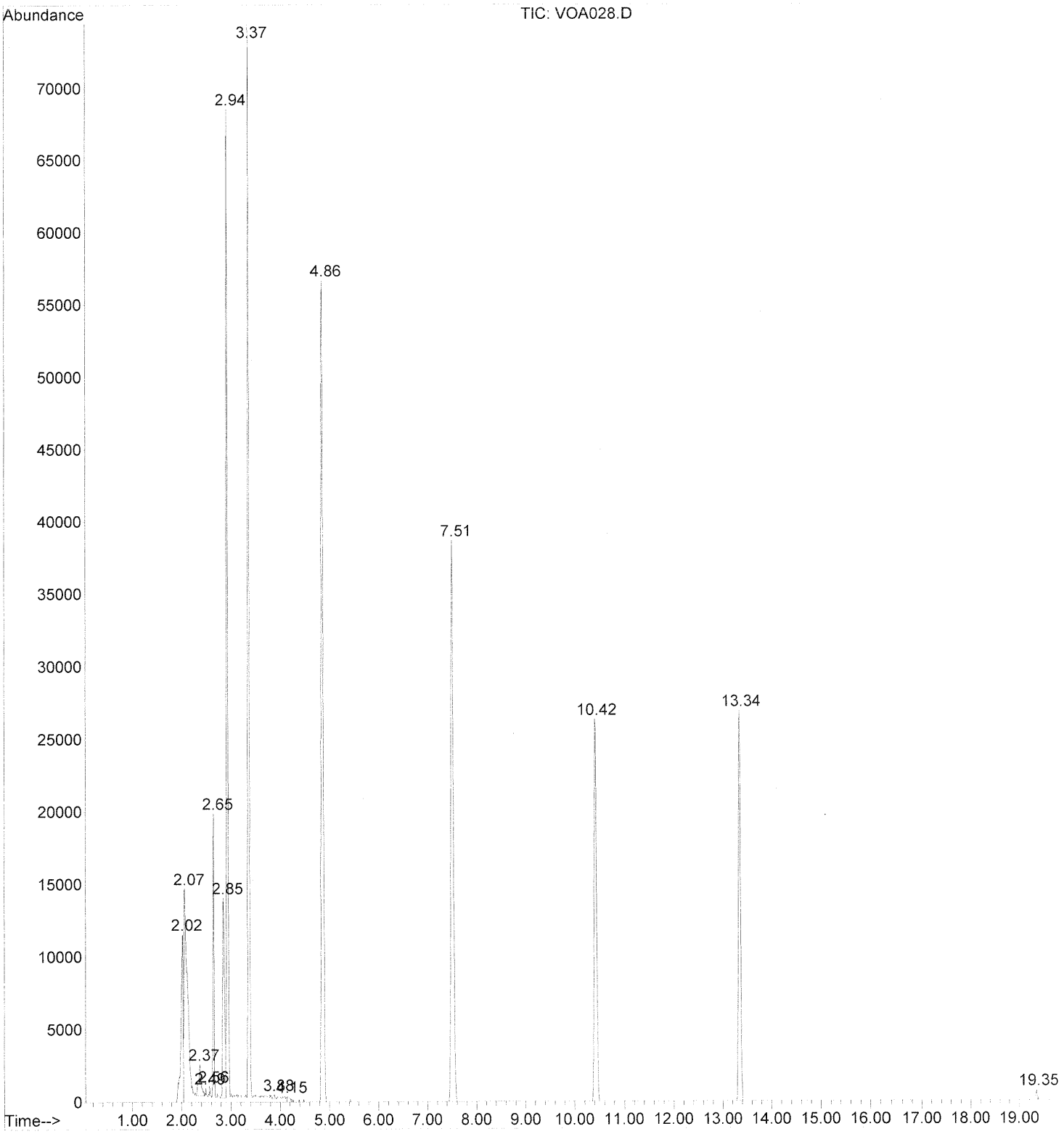
#	Compound	Ret Time	Concentration	Peak Type
1	Pentafluorobenzene	2.94	100.0 ppb	ISTD1
2	Dichlorofluoromethane	***** NOT FOUND *****		
3	Chloromethane	***** NOT FOUND *****		
4	Vinyl Chloride	***** NOT FOUND *****		
5	Bromomethane	***** NOT FOUND *****		
6	Chloroethane	***** NOT FOUND *****		
7	Trichlorofluoromethane	***** NOT FOUND *****		
8	Acetonitrile	***** NOT FOUND *****		
9	Diethyl ether	***** NOT FOUND *****		
10	1,1-Dichloroethene	***** NOT FOUND *****		
11	Acrylonitrile	***** NOT FOUND *****		
12	1,1,2-Trichlorotrifluoroethane	***** NOT FOUND *****		
13	Methylene Chloride	2.36	10.0 ppb	<
14	Iodomethane	***** NOT FOUND *****		
15	trans-1,2-Dichloroethene	***** NOT FOUND *****		
16	MTBE	***** NOT FOUND *****		
17	Allyl chloride	***** NOT FOUND *****		
18	1,1-Dichloroethane	***** NOT FOUND *****		
19	Carbon disulfide	***** NOT FOUND *****		
20	Chloroprene	***** NOT FOUND *****		
21	Methacrylonitrile	***** NOT FOUND *****		
22	cis-1,2-Dichloroethene	***** NOT FOUND *****		
23	Methyl acrylate	***** NOT FOUND *****		
24	Chloroform	***** NOT FOUND *****		
25	2,2-Dichloropropane	***** NOT FOUND *****		
26	Dibromofluoromethane-surrogate	2.84	109.6 ppb	110 %R
27	Isobutyl alcohol	***** NOT FOUND *****		
28	Tetrahydrofuran	***** NOT FOUND *****		
29	1,1,1-Trichloroethane	***** NOT FOUND *****		
30	1,2-Dichloroethane	***** NOT FOUND *****		
31	1,1-Dichloropropene	***** NOT FOUND *****		
32	Benzene	***** NOT FOUND *****		

33	Carbon Tetrachloride	***** NOT FOUND *****		
34	2-Chloroethanol	***** NOT FOUND *****		
35	1,4-Difluorobenzene	3.36	100.0 ppb	ISTD2
36	1,2-Dichloropropane	***** NOT FOUND *****		
37	Trichloroethene	***** NOT FOUND *****		
38	Dibromomethane	***** NOT FOUND *****		
39	Bromodichloromethane	***** NOT FOUND *****		
40	Methyl methacrylate	***** NOT FOUND *****		
41	1,4-Dioxane	***** NOT FOUND *****		
42	cis-1,3-Dichloropropene	***** NOT FOUND *****		
43	2-Nitropropane	***** NOT FOUND *****		
44	trans-1,3-Dichloropropene	***** NOT FOUND *****		
45	Toluene	***** NOT FOUND *****		
46	Toluene-d8-surrogate	4.86	103.9 ppb	104 % R
47	1,1,2-Trichloroethane	***** NOT FOUND *****		
48	1,3-Dichloropropane	***** NOT FOUND *****		
49	Ethylmethacrylate	***** NOT FOUND *****		
50	Dibromochloromethane	***** NOT FOUND *****		
51	1,2-Dibromomethane	***** NOT FOUND *****		
52	Tetrachloroethene	***** NOT FOUND *****		
53	Chlorobenzene	***** NOT FOUND *****		
54	Chlorobenzene-d5	7.52	100.0 ppb	ISTD3
55	1,1,1,2-Tetrachloroethane	***** NOT FOUND *****		
56	Ethylbenzene	***** NOT FOUND *****		
57	m/p-xylene	***** NOT FOUND *****		
58	Bromoform	***** NOT FOUND *****		
59	Styrene	***** NOT FOUND *****		
60	o-Xylene	***** NOT FOUND *****		
61	cis-1,4-Dichloro-2-butene	***** NOT FOUND *****		
62	trans-1,4-Dichloro-2-butene	***** NOT FOUND *****		82 % R
63	Bromofluorobenzene-surrogate	10.45	81.5 ppb	
64	1,1,2,2-Tetrachloroethane	***** NOT FOUND *****		
65	1,2,3-Trichloropropane	***** NOT FOUND *****		
66	Isopropylbenzene	***** NOT FOUND *****		
67	Bromobenzene	***** NOT FOUND *****		
68	2-Chlorotoluene	***** NOT FOUND *****		
69	n-Propylbenzene	***** NOT FOUND *****		
70	4-Chlorotoluene	***** NOT FOUND *****		
71	1,3,5-Trimethylbenzene	***** NOT FOUND *****		
72	Pentachloroethane	***** NOT FOUND *****		
73	tert-Butylbenzene	***** NOT FOUND *****		
74	1,2,4-Trimethylbenzene	***** NOT FOUND *****		
75	1,3-Dichlorobenzene	***** NOT FOUND *****		
76	1,4-Dichlorobenzene-d4	13.34	100.0 ppb	ISTD4
77	sec-Butylbenzene	***** NOT FOUND *****		
78	1,4-Dichlorobenzene	***** NOT FOUND *****		
79	p-iso-Propyltoluene	***** NOT FOUND *****		
80	1,2-Dichlorobenzene	***** NOT FOUND *****		
81	n-Butylbenzene	***** NOT FOUND *****		
82	1,2-Bromo-3-chloropropane	***** NOT FOUND *****		
83	Nitobenzene	***** NOT FOUND *****		
84	1,2,4-Trichlorobenzene	***** NOT FOUND *****		
85	Naphthalene	***** NOT FOUND *****		
86	1,2,3-Trichlorobenzene	***** NOT FOUND *****		
87	Hexachlorobutadiene	***** NOT FOUND *****		

END OF REPORT

Thu May 25 14:09:45 2006

File : C:\HPCHEM\1\DATA\052406\VOA028.D
Operator : dmt
Acquired : 25 May 2006 2:15 pm using AcqMethod VOA0506
Instrument : GCD Syste
Sample Name: SB-1 0-4' SOIL
Misc Info :
Vial Number: 1



Calibration Last Updated: Thu May 25 13:44:00 2006
Quantitation Report - Summary

Information from Data File:
File : C:\HPCHEM\1\DATA\052406\VOA028.D
Operator : dmt
Acquired : 25 May 2006 2:15 pm using AcqMethod VOA0506
Sample Name: SB-1 0-4' SOIL
Misc Info :
Vial Number: 1
CurrentMeth: C:\HPCHEM\1\METHODS\VOA0506R.M

Quantitation Settings:
Non-Reference Peak Window: 10.00 Percent
Correlation Window: 0.02 Minutes
Default Multiplier: 1.00
Default Sample Concentration: 0.00
Peak Type Decoding: m -> Manually re-integrated

Number of Compounds in Database: 87

#	Compound	Ret Time	Concentration	Peak Type
1	Pentafluorobenzene	2.94	100.0 ppb	ISTD1#
2	Dichlorofluoromethane	***** NOT FOUND *****		
3	Chloromethane	***** NOT FOUND *****		
4	Vinyl Chloride	***** NOT FOUND *****		
5	Bromomethane	***** NOT FOUND *****		
6	Chloroethane	***** NOT FOUND *****		
7	Trichlorofluoromethane	***** NOT FOUND *****		
8	Acetonitrile	***** NOT FOUND *****		
9	Diethyl ether	***** NOT FOUND *****		
10	1,1-Dichloroethene	~~2.43~~	~~3.0 ppb~~	~~#~~ m
11	Acrylonitrile	***** NOT FOUND *****		
12	1,1,2-Trichlorotrifluoroethane	***** NOT FOUND *****		
13	Methylene Chloride	2.38	6.0 ppb	#
14	Iodomethane	***** NOT FOUND *****		
15	trans-1,2-Dichloroethene	~~2.43~~	~~3.1 ppb~~	~~#~~ m
16	MTBE	***** NOT FOUND *****		
17	Allyl chloride	***** NOT FOUND *****		
18	1,1-Dichloroethane	***** NOT FOUND *****		
19	Carbon disulfide	***** NOT FOUND *****		
20	Chloroprene	***** NOT FOUND *****		
21	Methacrylonitrile	***** NOT FOUND *****		
22	cis-1,2-Dichloroethene	***** NOT FOUND *****		
23	Methyl acrylate	***** NOT FOUND *****		
24	Chloroform	***** NOT FOUND *****		
25	2,2-Dichloropropane	***** NOT FOUND *****		
26	Dibromofluoromethane-surrogate	2.85	79.1 ppb	79 %
27	Isobutyl alcohol	***** NOT FOUND *****		
28	Tetrahydrofuran	***** NOT FOUND *****		
29	1,1,1-Trichloroethane	***** NOT FOUND *****		
30	1,2-Dichloroethane	***** NOT FOUND *****		
31	1,1-Dichloropropene	***** NOT FOUND *****		
32	Benzene	***** NOT FOUND *****		

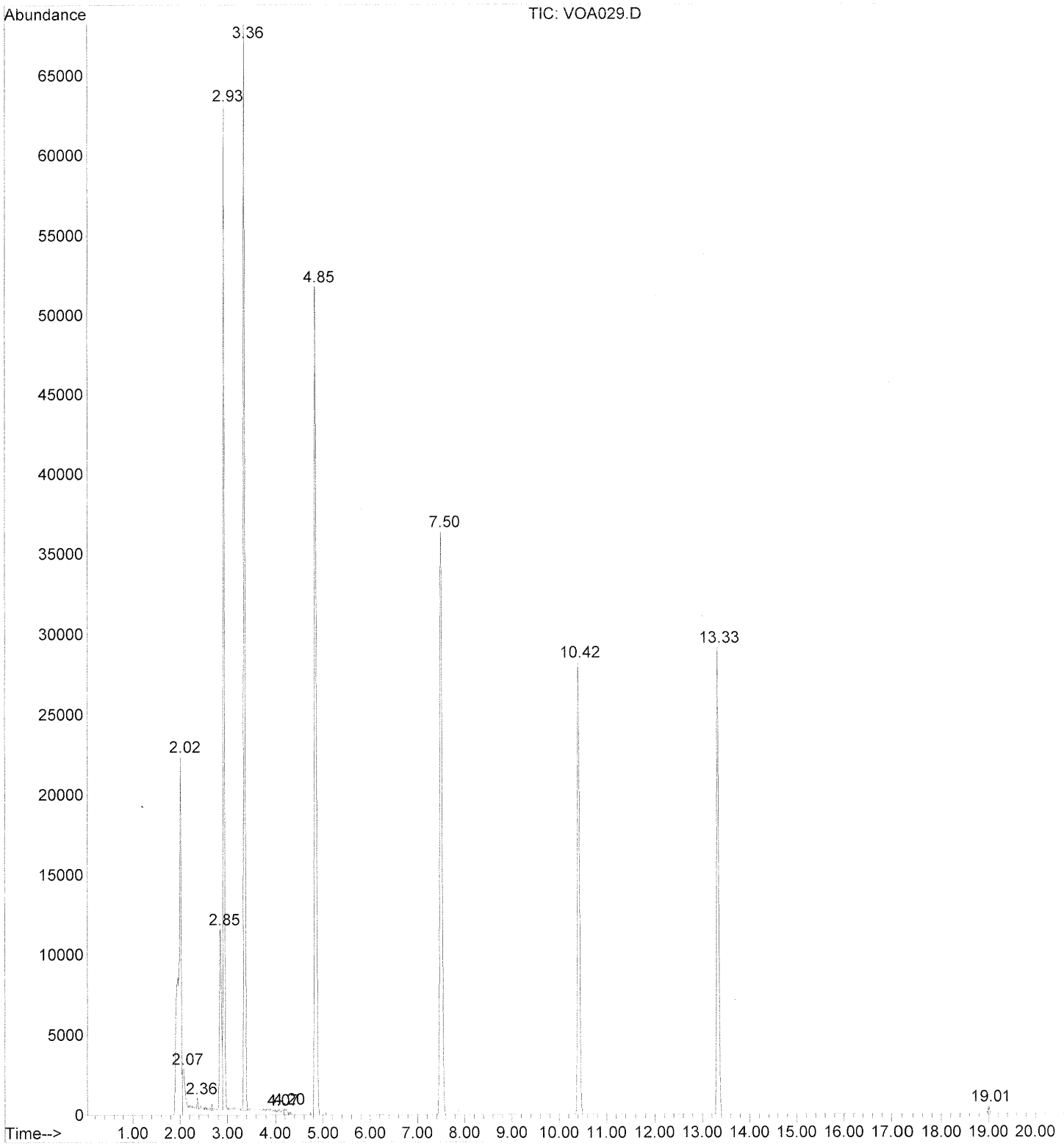
33	Carbon Tetrachloride	***** NOT FOUND *****		
34	2-Chloroethanol	***** NOT FOUND *****		
35	1,4-Difluorobenzene	3.37	100.0 ppb	ISTD2#
36	1,2-Dichloropropane	***** NOT FOUND *****		
37	Trichloroethene	***** NOT FOUND *****		
38	Dibromomethane	***** NOT FOUND *****		
39	Bromodichloromethane	***** NOT FOUND *****		
40	Methyl methacrylate	***** NOT FOUND *****		
41	1,4-Dioxane	***** NOT FOUND *****		
42	cis-1,3-Dichloropropene	***** NOT FOUND *****		
43	2-Nitropropane	***** NOT FOUND *****		
44	trans-1,3-Dichloropropene	***** NOT FOUND *****		
45	Toluene	***** NOT FOUND *****		
46	Toluene-d8-surrogate	4.86	97.0 ppb	97 % L
47	1,1,2-Trichloroethane	***** NOT FOUND *****		
48	1,3-Dichloropropane	***** NOT FOUND *****		
49	Ethylmethacrylate	***** NOT FOUND *****		
50	Dibromochloromethane	***** NOT FOUND *****		
51	1,2-Dibromomethane	***** NOT FOUND *****		
52	Tetrachloroethene	***** NOT FOUND *****		
53	Chlorobenzene	***** NOT FOUND *****		
54	Chlorobenzene-d5	7.51	100.0 ppb	ISTD3
55	1,1,1,2-Tetrachloroethane	***** NOT FOUND *****		
56	Ethylbenzene	***** NOT FOUND *****		
57	m/p-xylene	***** NOT FOUND *****		
58	Bromoform	***** NOT FOUND *****		
59	Styrene	***** NOT FOUND *****		
60	o-Xylene	***** NOT FOUND *****		
61	cis-1,4-Dichloro-2-butene	***** NOT FOUND *****		
62	trans-1,4-Dichloro-2-butene	***** NOT FOUND *****		
63	Bromofluorobenzene-surrogate	10.42	84.9 ppb	85 % L
64	1,1,2,2-Tetrachloroethane	***** NOT FOUND *****		
65	1,2,3-Trichloropropane	***** NOT FOUND *****		
66	Isopropylbenzene	***** NOT FOUND *****		
67	Bromobenzene	***** NOT FOUND *****		
68	2-Chlorotoluene	***** NOT FOUND *****		
69	n-Propylbenzene	***** NOT FOUND *****		
70	4-Chlorotoluene	***** NOT FOUND *****		
71	1,3,5-Trimethylbenzene	***** NOT FOUND *****		
72	Pentachloroethane	***** NOT FOUND *****		
73	tert-Butylbenzene	***** NOT FOUND *****		
74	1,2,4-Trimethylbenzene	***** NOT FOUND *****		
75	1,3-Dichlorobenzene	***** NOT FOUND *****		
76	1,4-Dichlorobenzene-d4	13.34	100.0 ppb	ISTD4
77	sec-Butylbenzene	***** NOT FOUND *****		
78	1,4-Dichlorobenzene	***** NOT FOUND *****		
79	p-iso-Propyltoluene	***** NOT FOUND *****		
80	1,2-Dichlorobenzene	***** NOT FOUND *****		
81	n-Butylbenzene	***** NOT FOUND *****		
82	1,2-Bromo-3-chloropropane	***** NOT FOUND *****		
83	Nitobenzene	***** NOT FOUND *****		
84	1,2,4-Trichlorobenzene	***** NOT FOUND *****		
85	Naphthalene	***** NOT FOUND *****		
86	1,2,3-Trichlorobenzene	***** NOT FOUND *****		
87	Hexachlorobutadiene	***** NOT FOUND *****		

END OF REPORT

Qualifiers Not Satisfied

Thu May 25 14:38:35 2006

File : C:\HPCHEM\1\DATA\052406\VOA029.D
Operator : dmt
Acquired : 25 May 2006 2:42 pm using AcqMethod VOA0506
Instrument : GCD Syste
Sample Name: SB-38 8-12' H2O
Misc Info :
Vial Number: 1



Calibration Last Updated: Thu May 25 13:44:00 2006
Quantitation Report - Summary

Information from Data File:
File : C:\HPCHEM\1\DATA\052406\VOA029.D
Operator : dmt
Acquired : 25 May 2006 2:42 pm using AcqMethod VOA0506
Sample Name: SB-38 8-12' H2O
Misc Info :
Vial Number: 1
CurrentMeth: C:\HPCHEM\1\METHODS\VOA0506R.M

Quantitation Settings:
Non-Reference Peak Window: 10.00 Percent
Correlation Window: 0.02 Minutes
Default Multiplier: 1.00
Default Sample Concentration: 0.00
Peak Type Decoding: m -> Manually re-integrated

Number of Compounds in Database: 87

#	Compound	Ret Time	Concentration	Peak Type
1	Pentafluorobenzene	2.93	100.0 ppb	ISTD1
2	Dichlorofluoromethane	***** NOT FOUND *****		
3	Chloromethane	***** NOT FOUND *****		
4	Vinyl Chloride	***** NOT FOUND *****		
5	Bromomethane	***** NOT FOUND *****		
6	Chloroethane	***** NOT FOUND *****		
7	Trichlorofluoromethane	***** NOT FOUND *****		
8	Acetonitrile	***** NOT FOUND *****		
9	Diethyl ether	***** NOT FOUND *****		
10	1,1-Dichloroethene	***** NOT FOUND *****		
11	Acrylonitrile	***** NOT FOUND *****		
12	1,1,2-Trichlorotrifluoroethane	***** NOT FOUND *****		
13	Methylene Chloride	2.37	3.3 ppb	#
14	Iodomethane	***** NOT FOUND *****		
15	trans-1,2-Dichloroethene	***** NOT FOUND *****		
16	MTBE	***** NOT FOUND *****		
17	Allyl chloride	***** NOT FOUND *****		
18	1,1-Dichloroethane	***** NOT FOUND *****		
19	Carbon disulfide	***** NOT FOUND *****		
20	Chloroprene	***** NOT FOUND *****		
21	Methacrylonitrile	***** NOT FOUND *****		
22	cis-1,2-Dichloroethene	***** NOT FOUND *****		
23	Methyl acrylate	***** NOT FOUND *****		
24	Chloroform	***** NOT FOUND *****		
25	2,2-Dichloropropane	***** NOT FOUND *****		
26	Dibromofluoromethane-surrogate	2.85	77.6 ppb	
27	Isobutyl alcohol	***** NOT FOUND *****		
28	Tetrahydrofuran	***** NOT FOUND *****		
29	1,1,1-Trichloroethane	***** NOT FOUND *****		
30	1,2-Dichloroethane	***** NOT FOUND *****		
31	1,1-Dichloropropene	***** NOT FOUND *****		
32	Benzene	***** NOT FOUND *****		

79 1-K

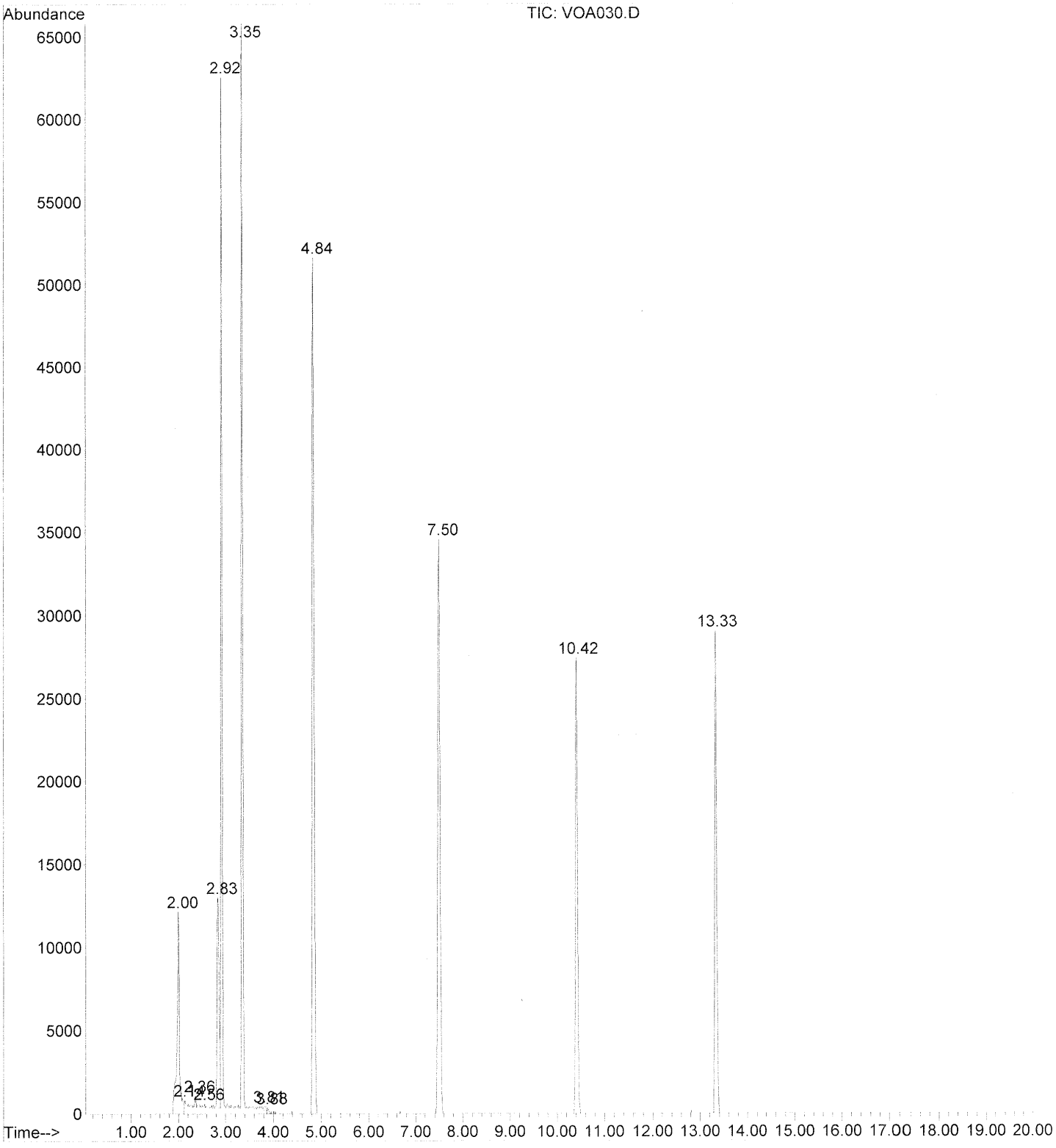
33	Carbon Tetrachloride	***** NOT FOUND *****		
34	2-Chloroethanol	***** NOT FOUND *****		
35	1,4-Difluorobenzene	3.36	100.0 ppb	ISTD2
36	1,2-Dichloropropane	***** NOT FOUND *****		
37	Trichloroethene	***** NOT FOUND *****		
38	Dibromomethane	***** NOT FOUND *****		
39	Bromodichloromethane	***** NOT FOUND *****		
40	Methyl methacrylate	***** NOT FOUND *****		
41	1,4-Dioxane	***** NOT FOUND *****		
42	cis-1,3-Dichloropropene	***** NOT FOUND *****		
43	2-Nitropropane	***** NOT FOUND *****		
44	trans-1,3-Dichloropropene	***** NOT FOUND *****		
45	Toluene	***** NOT FOUND *****		
46	Toluene-d8-surrogate	4.85	100.5 ppb	101 %R
47	1,1,2-Trichloroethane	***** NOT FOUND *****		
48	1,3-Dichloropropane	***** NOT FOUND *****		
49	Ethylmethacrylate	***** NOT FOUND *****		
50	Dibromochloromethane	***** NOT FOUND *****		
51	1,2-Dibromomethane	***** NOT FOUND *****		
52	Tetrachloroethene	***** NOT FOUND *****		
53	Chlorobenzene	***** NOT FOUND *****		
54	Chlorobenzene-d5	7.51	100.0 ppb	ISTD3
55	1,1,1,2-Tetrachloroethane	***** NOT FOUND *****		
56	Ethylbenzene	***** NOT FOUND *****		
57	m/p-xylene	***** NOT FOUND *****		
58	Bromoform	***** NOT FOUND *****		
59	Styrene	***** NOT FOUND *****		
60	o-Xylene	***** NOT FOUND *****		
61	cis-1,4-Dichloro-2-butene	***** NOT FOUND *****		
62	trans-1,4-Dichloro-2-butene	***** NOT FOUND *****		
63	Bromofluorobenzene-surrogate	10.42	86.0 ppb	86 %R
64	1,1,2,2-Tetrachloroethane	***** NOT FOUND *****		
65	1,2,3-Trichloropropane	***** NOT FOUND *****		
66	Isopropylbenzene	***** NOT FOUND *****		
67	Bromobenzene	***** NOT FOUND *****		
68	2-Chlorotoluene	***** NOT FOUND *****		
69	n-Propylbenzene	***** NOT FOUND *****		
70	4-Chlorotoluene	***** NOT FOUND *****		
71	1,3,5-Trimethylbenzene	***** NOT FOUND *****		
72	Pentachloroethane	***** NOT FOUND *****		
73	tert-Butylbenzene	***** NOT FOUND *****		
74	1,2,4-Trimethylbenzene	***** NOT FOUND *****		
75	1,3-Dichlorobenzene	***** NOT FOUND *****		
76	1,4-Dichlorobenzene-d4	13.33	100.0 ppb	ISTD4
77	sec-Butylbenzene	***** NOT FOUND *****		
78	1,4-Dichlorobenzene	***** NOT FOUND *****		
79	p-iso-Propyltoluene	***** NOT FOUND *****		
80	1,2-Dichlorobenzene	***** NOT FOUND *****		
81	n-Butylbenzene	***** NOT FOUND *****		
82	1,2-Bromo-3-chloropropane	***** NOT FOUND *****		
83	Nitobenzene	***** NOT FOUND *****		
84	1,2,4-Trichlorobenzene	***** NOT FOUND *****		
85	Naphthalene	***** NOT FOUND *****		
86	1,2,3-Trichlorobenzene	***** NOT FOUND *****		
87	Hexachlorobutadiene	***** NOT FOUND *****		

END OF REPORT

Qualifiers Not Satisfied

Thu May 25 15:11:26 2006

File : C:\HPCHEM\1\DATA\052406\VOA030.D
Operator : dmt
Acquired : 25 May 2006 3:24 pm using AcqMethod VOA0506
Instrument : GCD Syste
Sample Name: SB-37 7.3-11.7' H2O
Misc Info :
Vial Number: 1



Calibration Last Updated: Thu May 25 13:44:00 2006
Quantitation Report - Summary

Information from Data File:
File : C:\HPCHEM\1\DATA\052406\VOA030.D
Operator : dmt
Acquired : 25 May 2006 3:24 pm using AcqMethod VOA0506
Sample Name: SB-37 7.3-11.7' H2O
Misc Info :
Vial Number: 1
CurrentMeth: C:\HPCHEM\1\METHODS\VOA0506R.M

Quantitation Settings:
Non-Reference Peak Window: 10.00 Percent
Correlation Window: 0.02 Minutes
Default Multiplier: 1.00
Default Sample Concentration: 0.00
Peak Type Decoding: m -> Manually re-integrated

Number of Compounds in Database: 87

#	Compound	Ret Time	Concentration	Peak Type
1	Pentafluorobenzene	2.92	100.0 ppb	ISTD1#
2	Dichlorofluoromethane	***** NOT FOUND *****		
3	Chloromethane	***** NOT FOUND *****		
4	Vinyl Chloride	***** NOT FOUND *****		
5	Bromomethane	***** NOT FOUND *****		
6	Chloroethane	***** NOT FOUND *****		
7	Trichlorofluoromethane	***** NOT FOUND *****		
8	Acetonitrile	***** NOT FOUND *****		
9	Diethyl ether	***** NOT FOUND *****		
10	1,1-Dichloroethene	***** NOT FOUND *****		
11	Acrylonitrile	***** NOT FOUND *****		
12	1,1,2-Trichlorotrifluoroethane	***** NOT FOUND *****		
13	Methylene Chloride	2.36	4.0 ppb	#
14	Iodomethane	***** NOT FOUND *****		
15	trans-1,2-Dichloroethene	***** NOT FOUND *****		
16	MTBE	***** NOT FOUND *****		
17	Allyl chloride	***** NOT FOUND *****		
18	1,1-Dichloroethane	***** NOT FOUND *****		
19	Carbon disulfide	***** NOT FOUND *****		
20	Chloroprene	***** NOT FOUND *****		
21	Methacrylonitrile	***** NOT FOUND *****		
22	cis-1,2-Dichloroethene	***** NOT FOUND *****		
23	Methyl acrylate	***** NOT FOUND *****		
24	Chloroform	***** NOT FOUND *****		
25	2,2-Dichloropropane	***** NOT FOUND *****		
26	Dibromofluoromethane-surrogate	2.83	83.9 ppb	
27	Isobutyl alcohol	***** NOT FOUND *****		
28	Tetrahydrofuran	***** NOT FOUND *****		
29	1,1,1-Trichloroethane	***** NOT FOUND *****		
30	1,2-Dichloroethane	***** NOT FOUND *****		
31	1,1-Dichloropropene	***** NOT FOUND *****		
32	Benzene	***** NOT FOUND *****		

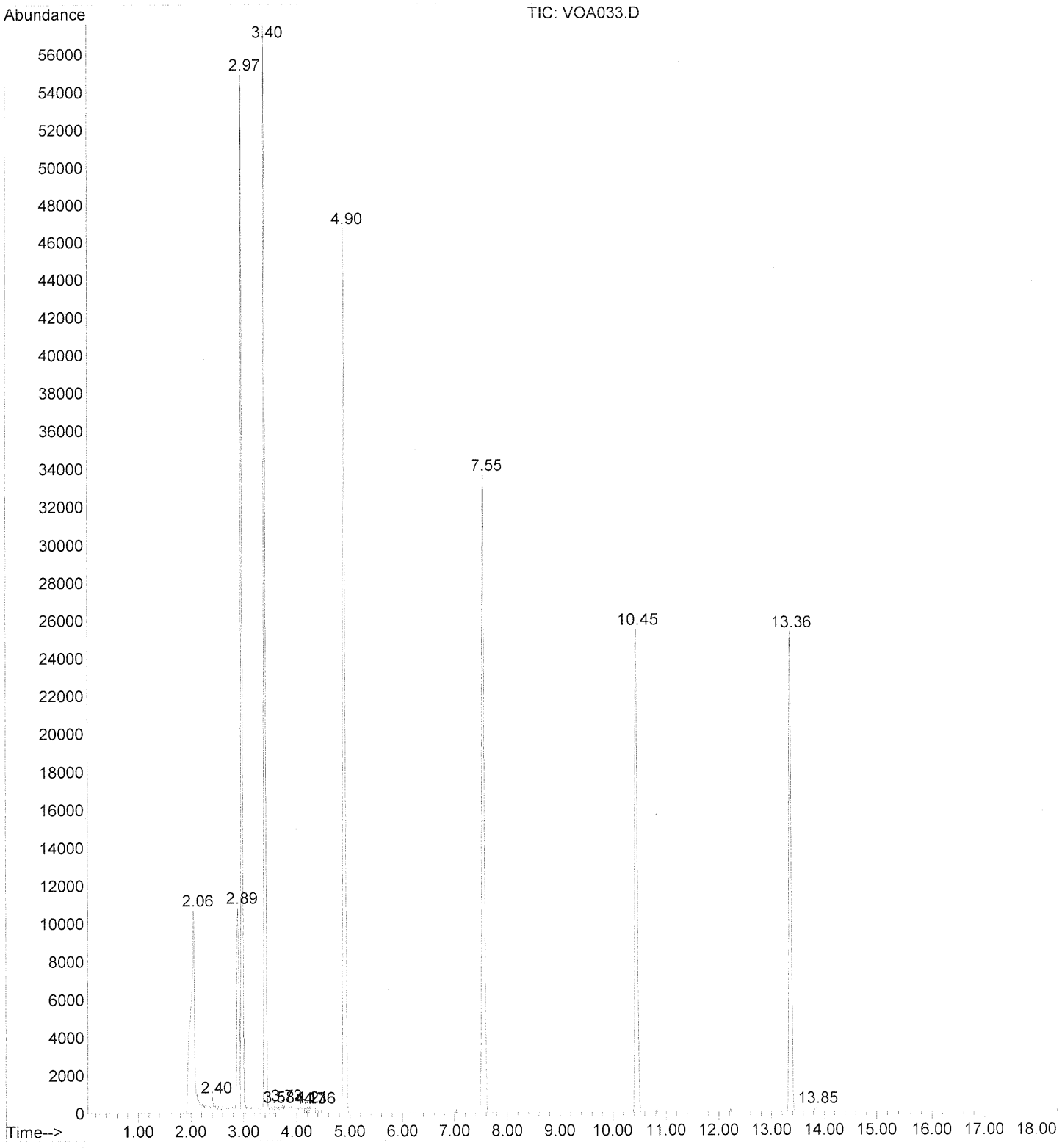
33	Carbon Tetrachloride	***** NOT FOUND *****		
34	2-Chloroethanol	***** NOT FOUND *****		
35	1,4-Difluorobenzene	3.35	100.0 ppb	ISTD2#
36	1,2-Dichloropropane	***** NOT FOUND *****		
37	Trichloroethene	***** NOT FOUND *****		
38	Dibromomethane	***** NOT FOUND *****		
39	Bromodichloromethane	***** NOT FOUND *****		
40	Methyl methacrylate	***** NOT FOUND *****		
41	1,4-Dioxane	***** NOT FOUND *****		
42	cis-1,3-Dichloropropene	***** NOT FOUND *****		
43	2-Nitropropane	***** NOT FOUND *****		
44	trans-1,3-Dichloropropene	***** NOT FOUND *****		
45	Toluene	***** NOT FOUND *****		98 1/2
46	Toluene-d8-surrogate	4.84	98.2 ppb	
47	1,1,2-Trichloroethane	***** NOT FOUND *****		
48	1,3-Dichloropropane	***** NOT FOUND *****		
49	Ethylmethacrylate	***** NOT FOUND *****		
50	Dibromochloromethane	***** NOT FOUND *****		
51	1,2-Dibromomethane	***** NOT FOUND *****		
52	Tetrachloroethene	***** NOT FOUND *****		
53	Chlorobenzene	***** NOT FOUND *****		
54	Chlorobenzene-d5	7.50	100.0 ppb	ISTD3
55	1,1,1,2-Tetrachloroethane	***** NOT FOUND *****		
56	Ethylbenzene	***** NOT FOUND *****		
57	m/p-xylene	***** NOT FOUND *****		
58	Bromoform	***** NOT FOUND *****		
59	Styrene	***** NOT FOUND *****		
60	o-Xylene	***** NOT FOUND *****		
61	cis-1,4-Dichloro-2-butene	***** NOT FOUND *****		
62	trans-1,4-Dichloro-2-butene	***** NOT FOUND *****		
63	Bromofluorobenzene-surrogate	10.42	91.9 ppb	92 1/2
64	1,1,2,2-Tetrachloroethane	***** NOT FOUND *****		
65	1,2,3-Trichloropropane	***** NOT FOUND *****		
66	Isopropylbenzene	***** NOT FOUND *****		
67	Bromobenzene	***** NOT FOUND *****		
68	2-Chlorotoluene	***** NOT FOUND *****		
69	n-Propylbenzene	***** NOT FOUND *****		
70	4-Chlorotoluene	***** NOT FOUND *****		
71	1,3,5-Trimethylbenzene	***** NOT FOUND *****		
72	Pentachloroethane	***** NOT FOUND *****		
73	tert-Butylbenzene	***** NOT FOUND *****		
74	1,2,4-Trimethylbenzene	***** NOT FOUND *****		
75	1,3-Dichlorobenzene	***** NOT FOUND *****		
76	1,4-Dichlorobenzene-d4	13.33	100.0 ppb	ISTD4
77	sec-Butylbenzene	***** NOT FOUND *****		
78	1,4-Dichlorobenzene	***** NOT FOUND *****		
79	p-iso-Propyltoluene	***** NOT FOUND *****		
80	1,2-Dichlorobenzene	***** NOT FOUND *****		
81	n-Butylbenzene	***** NOT FOUND *****		
82	1,2-Bromo-3-chloropropane	***** NOT FOUND *****		
83	Nitobenzene	***** NOT FOUND *****		
84	1,2,4-Trichlorobenzene	***** NOT FOUND *****		
85	Naphthalene	***** NOT FOUND *****		
86	1,2,3-Trichlorobenzene	***** NOT FOUND *****		
87	Hexachlorobutadiene	***** NOT FOUND *****		

END OF REPORT

Qualifiers Not Satisfied

Thu May 25 15:46:14 2006

File : C:\HPCHEM\1\DATA\052406\VOA033.D
Operator : dmt
Acquired : 26 May 2006 8:45 am using AcqMethod VOA0506
Instrument : GCD Syste
Sample Name: MW-4 H2O
Misc Info :
Vial Number: 1



Calibration Last Updated: Fri May 26 08:11:46 2006
Quantitation Report - Summary

Information from Data File:
File : C:\HPCHEM\1\DATA\052406\VOA033.D
Operator : dmt
Acquired : 26 May 2006 8:45 am using AcqMethod VOA0506
Sample Name: MW-4 H2O
Misc Info :
Vial Number: 1
CurrentMeth: C:\HPCHEM\1\METHODS\VOA0506R.M

Quantitation Settings:
Non-Reference Peak Window: 10.00 Percent
Correlation Window: 0.02 Minutes
Default Multiplier: 1.00
Default Sample Concentration: 0.00
Peak Type Decoding: m -> Manually re-integrated

Number of Compounds in Database: 87

#	Compound	Ret Time	Concentration	Peak Type
1	Pentafluorobenzene	2.97	100.0 ppb	ISTD1
2	Dichlorofluoromethane	***** NOT FOUND *****		
3	Chloromethane	***** NOT FOUND *****		
4	Vinyl Chloride	***** NOT FOUND *****		
5	Bromomethane	***** NOT FOUND *****		
6	Chloroethane	***** NOT FOUND *****		
7	Trichlorofluoromethane	***** NOT FOUND *****		
8	Acetonitrile	***** NOT FOUND *****		
9	Diethyl ether	***** NOT FOUND *****		
10	1,1-Dichloroethene	***** NOT FOUND *****		
11	Acrylonitrile	***** NOT FOUND *****		
12	1,1,2-Trichlorotrifluoroethane	***** NOT FOUND *****		
13	Methylene Chloride	2.41	5.1 ppb	#
14	Iodomethane	***** NOT FOUND *****		
15	trans-1,2-Dichloroethene	***** NOT FOUND *****		
16	MTBE	***** NOT FOUND *****		
17	Allyl chloride	***** NOT FOUND *****		
18	1,1-Dichloroethane	***** NOT FOUND *****		
19	Carbon disulfide	***** NOT FOUND *****		
20	Chloroprene	***** NOT FOUND *****		
21	Methacrylonitrile	***** NOT FOUND *****		
22	cis-1,2-Dichloroethene	***** NOT FOUND *****		
23	Methyl acrylate	***** NOT FOUND *****		
24	Chloroform	***** NOT FOUND *****		
25	2,2-Dichloropropane	***** NOT FOUND *****		
26	Dibromofluoromethane-surrogate	2.89	95.8 ppb	
27	Isobutyl alcohol	***** NOT FOUND *****		
28	Tetrahydrofuran	***** NOT FOUND *****		
29	1,1,1-Trichloroethane	***** NOT FOUND *****		
30	1,2-Dichloroethane	***** NOT FOUND *****		
31	1,1-Dichloropropene	***** NOT FOUND *****		
32	Benzene	***** NOT FOUND *****		

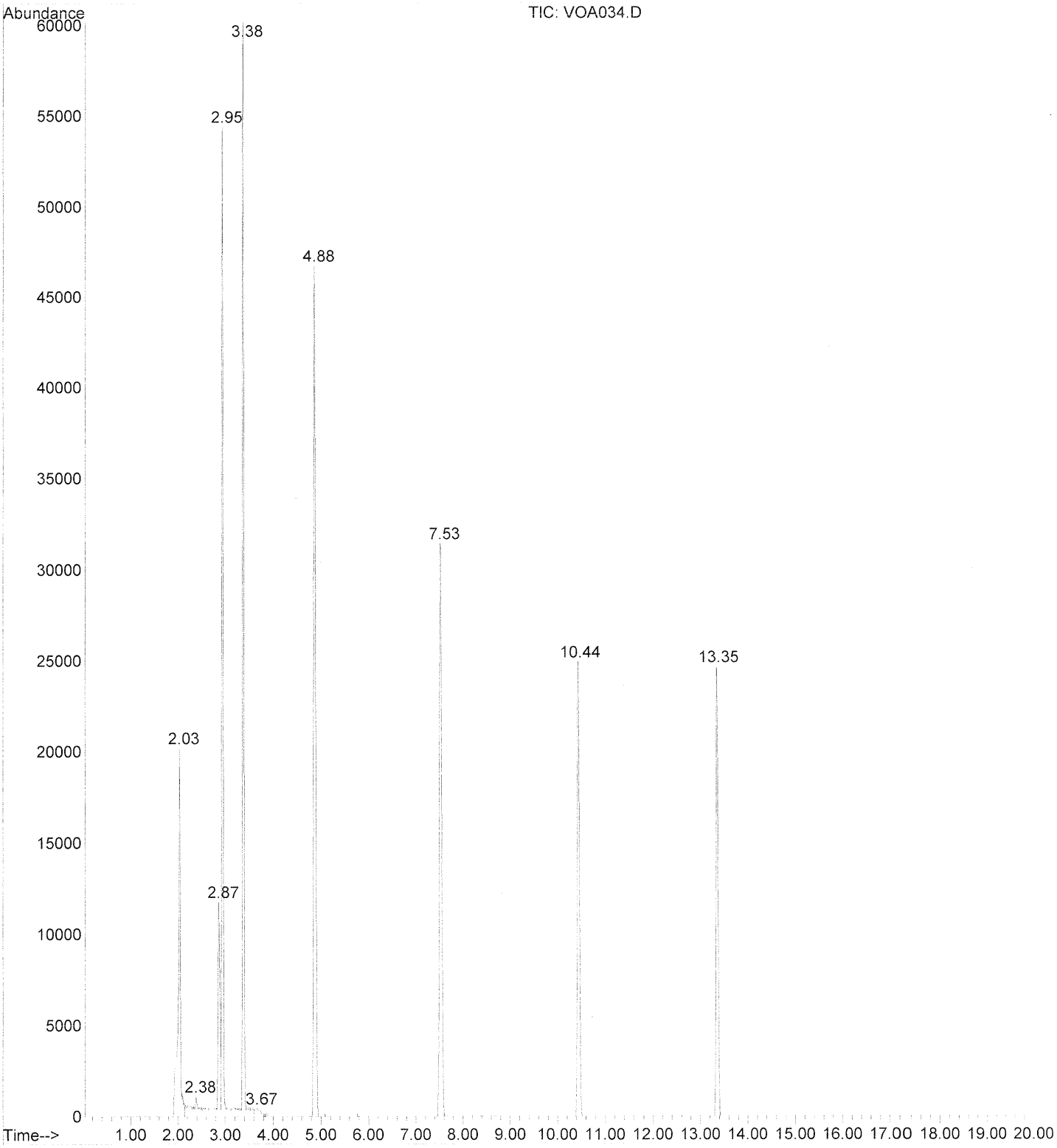
33	Carbon Tetrachloride	***** NOT FOUND *****		
34	2-Chloroethanol	***** NOT FOUND *****		
35	1,4-Difluorobenzene	3.40	100.0 ppb	ISTD2
36	1,2-Dichloropropane	***** NOT FOUND *****		
37	Trichloroethene	***** NOT FOUND *****		
38	Dibromomethane	***** NOT FOUND *****		
39	Bromodichloromethane	***** NOT FOUND *****		
40	Methyl methacrylate	***** NOT FOUND *****		
41	1,4-Dioxane	***** NOT FOUND *****		
42	cis-1,3-Dichloropropene	***** NOT FOUND *****		
43	2-Nitropropane	***** NOT FOUND *****		
44	trans-1,3-Dichloropropene	***** NOT FOUND *****		
45	Toluene	***** NOT FOUND *****		
46	Toluene-d8-surrogate	4.90	112.3 ppb	112 %
47	1,1,2-Trichloroethane	***** NOT FOUND *****		
48	1,3-Dichloropropane	***** NOT FOUND *****		
49	Ethylmethacrylate	***** NOT FOUND *****		
50	Dibromochloromethane	***** NOT FOUND *****		
51	1,2-Dibromomethane	***** NOT FOUND *****		
52	Tetrachloroethene	***** NOT FOUND *****		
53	Chlorobenzene	***** NOT FOUND *****		
54	Chlorobenzene-d5	7.55	100.0 ppb	ISTD3
55	1,1,1,2-Tetrachloroethane	***** NOT FOUND *****		
56	Ethylbenzene	***** NOT FOUND *****		
57	m/p-xylene	***** NOT FOUND *****		
58	Bromoform	***** NOT FOUND *****		
59	Styrene	***** NOT FOUND *****		
60	o-Xylene	***** NOT FOUND *****		
61	cis-1,4-Dichloro-2-butene	***** NOT FOUND *****		
62	trans-1,4-Dichloro-2-butene	***** NOT FOUND *****		
63	Bromofluorobenzene-surrogate	10.45	98.4 ppb	98 %
64	1,1,2,2-Tetrachloroethane	***** NOT FOUND *****		
65	1,2,3-Trichloropropane	***** NOT FOUND *****		
66	Isopropylbenzene	***** NOT FOUND *****		
67	Bromobenzene	***** NOT FOUND *****		
68	2-Chlorotoluene	***** NOT FOUND *****		
69	n-Propylbenzene	***** NOT FOUND *****		
70	4-Chlorotoluene	***** NOT FOUND *****		
71	1,3,5-Trimethylbenzene	***** NOT FOUND *****		
72	Pentachloroethane	***** NOT FOUND *****		
73	tert-Butylbenzene	***** NOT FOUND *****		
74	1,2,4-Trimethylbenzene	***** NOT FOUND *****		
75	1,3-Dichlorobenzene	***** NOT FOUND *****		
76	1,4-Dichlorobenzene-d4	13.36	100.0 ppb	ISTD4
77	sec-Butylbenzene	***** NOT FOUND *****		
78	1,4-Dichlorobenzene	***** NOT FOUND *****		
79	p-iso-Propyltoluene	***** NOT FOUND *****		
80	1,2-Dichlorobenzene	***** NOT FOUND *****		
81	n-Butylbenzene	***** NOT FOUND *****		
82	1,2-Bromo-3-chloropropane	***** NOT FOUND *****		
83	Nitobenzene	***** NOT FOUND *****		
84	1,2,4-Trichlorobenzene	***** NOT FOUND *****		
85	Naphthalene	***** NOT FOUND *****		
86	1,2,3-Trichlorobenzene	***** NOT FOUND *****		
87	Hexachlorobutadiene	***** NOT FOUND *****		

END OF REPORT

Qualifiers Not Satisfied

Fri May 26 09:04:23 2006

File : C:\HPCHEM\1\DATA\052406\VOA034.D
Operator : dmt
Acquired : 26 May 2006 9:28 am using AcqMethod VOA0506
Instrument : GCD Syste
Sample Name: SB-40 1.5-5.5' H2O
Misc Info :
Vial Number: 1



Calibration Last Updated: Fri May 26 08:11:46 2006
Quantitation Report - Summary

Information from Data File:
File : C:\HPCHEM\1\DATA\052406\VOA034.D
Operator : dmt
Acquired : 26 May 2006 9:28 am using AcqMethod VOA0506
Sample Name: SB-40 1.5-5.5' H2O
Misc Info :
Vial Number: 1
CurrentMeth: C:\HPCHEM\1\METHODS\VOA0506R.M

Quantitation Settings:
Non-Reference Peak Window: 10.00 Percent
Correlation Window: 0.02 Minutes
Default Multiplier: 1.00
Default Sample Concentration: 0.00
Peak Type Decoding: m -> Manually re-integrated

Number of Compounds in Database: 87

#	Compound	Ret Time	Concentration	Peak Type
1	Pentafluorobenzene	2.95	100.0 ppb	ISTD1
2	Dichlorofluoromethane	***** NOT FOUND *****		
3	Chloromethane	***** NOT FOUND *****		
4	Vinyl Chloride	***** NOT FOUND *****		
5	Bromomethane	***** NOT FOUND *****		
6	Chloroethane	***** NOT FOUND *****		
7	Trichlorofluoromethane	***** NOT FOUND *****		
8	Acetonitrile	***** NOT FOUND *****		
9	Diethyl ether	***** NOT FOUND *****		
10	1,1-Dichloroethene	***** NOT FOUND *****		
11	Acrylonitrile	***** NOT FOUND *****		
12	1,1,2-Trichlorotrifluoroethane	***** NOT FOUND *****		
13	Methylene Chloride	2.38	5.5 ppb	#
14	Iodomethane	***** NOT FOUND *****		
15	trans-1,2-Dichloroethene	***** NOT FOUND *****		
16	MTBE	***** NOT FOUND *****		
17	Allyl chloride	***** NOT FOUND *****		
18	1,1-Dichloroethane	***** NOT FOUND *****		
19	Carbon disulfide	***** NOT FOUND *****		
20	Chloroprene	***** NOT FOUND *****		
21	Methacrylonitrile	***** NOT FOUND *****		
22	cis-1,2-Dichloroethene	***** NOT FOUND *****		
23	Methyl acrylate	***** NOT FOUND *****		
24	Chloroform	***** NOT FOUND *****		
25	2,2-Dichloropropane	***** NOT FOUND *****		
26	Dibromofluoromethane-surrogate	2.87	90.1 ppb	
27	Isobutyl alcohol	***** NOT FOUND *****		
28	Tetrahydrofuran	***** NOT FOUND *****		
29	1,1,1-Trichloroethane	***** NOT FOUND *****		
30	1,2-Dichloroethane	***** NOT FOUND *****		
31	1,1-Dichloropropene	***** NOT FOUND *****		
32	Benzene	***** NOT FOUND *****		

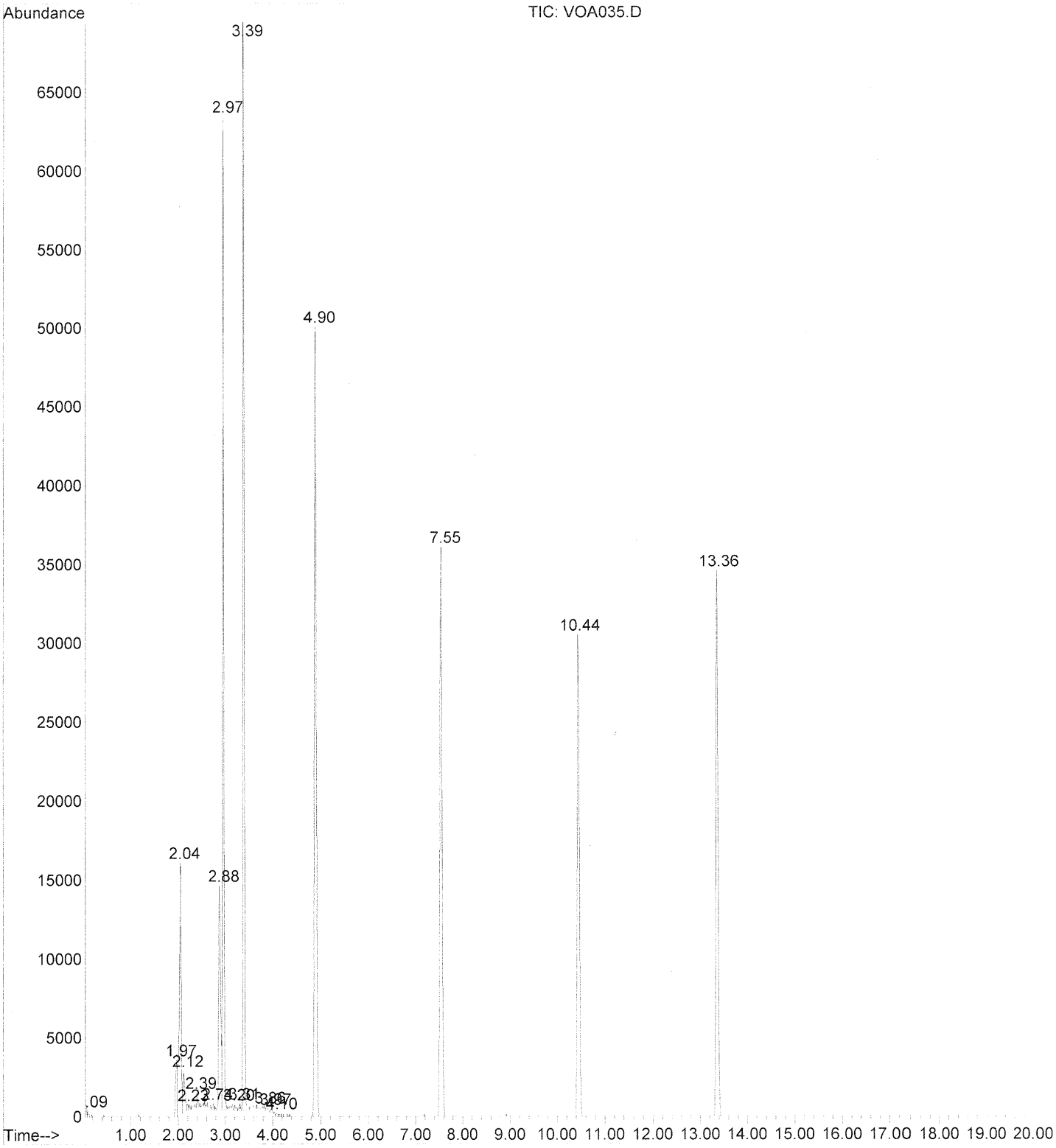
33	Carbon Tetrachloride	***** NOT FOUND *****		
34	2-Chloroethanol	***** NOT FOUND *****		
35	1,4-Difluorobenzene	3.38	100.0 ppb	ISTD2
36	1,2-Dichloropropane	***** NOT FOUND *****		
37	Trichloroethene	***** NOT FOUND *****		
38	Dibromomethane	***** NOT FOUND *****		
39	Bromodichloromethane	***** NOT FOUND *****		
40	Methyl methacrylate	***** NOT FOUND *****		
41	1,4-Dioxane	***** NOT FOUND *****		
42	cis-1,3-Dichloropropene	***** NOT FOUND *****		
43	2-Nitropropane	***** NOT FOUND *****		
44	trans-1,3-Dichloropropene	***** NOT FOUND *****		
45	Toluene	***** NOT FOUND *****		
46	Toluene-d8-surrogate	4.88	100.1 ppb	100 %R
47	1,1,2-Trichloroethane	***** NOT FOUND *****		
48	1,3-Dichloropropane	***** NOT FOUND *****		
49	Ethylmethacrylate	***** NOT FOUND *****		
50	Dibromochloromethane	***** NOT FOUND *****		
51	1,2-Dibromomethane	***** NOT FOUND *****		
52	Tetrachloroethene	***** NOT FOUND *****		
53	Chlorobenzene	***** NOT FOUND *****		
54	Chlorobenzene-d5	7.54	100.0 ppb	ISTD3
55	1,1,1,2-Tetrachloroethane	***** NOT FOUND *****		
56	Ethylbenzene	***** NOT FOUND *****		
57	m/p-xylene	***** NOT FOUND *****		
58	Bromoform	***** NOT FOUND *****		
59	Styrene	***** NOT FOUND *****		
60	o-Xylene	***** NOT FOUND *****		
61	cis-1,4-Dichloro-2-butene	***** NOT FOUND *****		
62	trans-1,4-Dichloro-2-butene	***** NOT FOUND *****		
63	Bromofluorobenzene-surrogate	10.44	95.2 ppb	95 %R
64	1,1,2,2-Tetrachloroethane	***** NOT FOUND *****		
65	1,2,3-Trichloropropane	***** NOT FOUND *****		
66	Isopropylbenzene	***** NOT FOUND *****		
67	Bromobenzene	***** NOT FOUND *****		
68	2-Chlorotoluene	***** NOT FOUND *****		
69	n-Propylbenzene	***** NOT FOUND *****		
70	4-Chlorotoluene	***** NOT FOUND *****		
71	1,3,5-Trimethylbenzene	***** NOT FOUND *****		
72	Pentachloroethane	***** NOT FOUND *****		
73	tert-Butylbenzene	***** NOT FOUND *****		
74	1,2,4-Trimethylbenzene	***** NOT FOUND *****		
75	1,3-Dichlorobenzene	***** NOT FOUND *****		
76	1,4-Dichlorobenzene-d4	13.35	100.0 ppb	ISTD4
77	sec-Butylbenzene	***** NOT FOUND *****		
78	1,4-Dichlorobenzene	***** NOT FOUND *****		
79	p-iso-Propyltoluene	***** NOT FOUND *****		
80	1,2-Dichlorobenzene	***** NOT FOUND *****		
81	n-Butylbenzene	***** NOT FOUND *****		
82	1,2-Bromo-3-chloropropane	***** NOT FOUND *****		
83	Nitobenzene	***** NOT FOUND *****		
84	1,2,4-Trichlorobenzene	***** NOT FOUND *****		
85	Naphthalene	***** NOT FOUND *****		
86	1,2,3-Trichlorobenzene	***** NOT FOUND *****		
87	Hexachlorobutadiene	***** NOT FOUND *****		

END OF REPORT

Qualifiers Not Satisfied

Fri May 26 09:49:28 2006

File : C:\HPCHEM\1\DATA\052406\VOA035.D
Operator : dmt
Acquired : 26 May 2006 12:03 pm using AcqMethod VOA0506
Instrument : GCD Syste
Sample Name: MW-5 4-11.6' H2O
Misc Info :
Vial Number: 1



Calibration Last Updated: Fri May 26 08:11:46 2006
Quantitation Report - Summary

Information from Data File:
File : C:\HPCHEM\1\DATA\052406\VOA035.D
Operator : dmt
Acquired : 26 May 2006 12:03 pm using AcqMethod VOA0506
Sample Name: MW-5 4-11.6' H2O
Misc Info :
Vial Number: 1
CurrentMeth: C:\HPCHEM\1\METHODS\VOA0506R.M

Quantitation Settings:
Non-Reference Peak Window: 10.00 Percent
Correlation Window: 0.02 Minutes
Default Multiplier: 1.00
Default Sample Concentration: 0.00
Peak Type Decoding: m -> Manually re-integrated

Number of Compounds in Database: 87

#	Compound	Ret Time	Concentration	Peak Type
1	Pentafluorobenzene	2.97	100.0 ppb	ISTD1#
2	Dichlorofluoromethane	***** NOT FOUND *****		
3	Chloromethane	***** NOT FOUND *****		
4	Vinyl Chloride	***** NOT FOUND *****		
5	Bromomethane	***** NOT FOUND *****		
6	Chloroethane	***** NOT FOUND *****		
7	Trichlorofluoromethane	***** NOT FOUND *****		
8	Acetonitrile	***** NOT FOUND *****		
9	Diethyl ether	***** NOT FOUND *****		
10	1,1-Dichloroethene	***** NOT FOUND *****		
11	Acrylonitrile	***** NOT FOUND *****		
12	1,1,2-Trichlorotrifluoroethane	***** NOT FOUND *****		
13	Methylene Chloride	2.40	3.7 ppb	#
14	Iodomethane	***** NOT FOUND *****		
15	trans-1,2-Dichloroethene	***** NOT FOUND *****		
16	MTBE	***** NOT FOUND *****		
17	Allyl chloride	***** NOT FOUND *****		
18	1,1-Dichloroethane	***** NOT FOUND *****		
19	Carbon disulfide	***** NOT FOUND *****		
20	Chloroprene	***** NOT FOUND *****		
21	Methacrylonitrile	***** NOT FOUND *****		
22	cis-1,2-Dichloroethene	***** NOT FOUND *****		
23	Methyl acrylate	***** NOT FOUND *****		
24	Chloroform	***** NOT FOUND *****		
25	2,2-Dichloropropane	***** NOT FOUND *****		
26	Dibromofluoromethane-surrogate	2.88	87.1 ppb	
27	Isobutyl alcohol	***** NOT FOUND *****		
28	Tetrahydrofuran	***** NOT FOUND *****		
29	1,1,1-Trichloroethane	***** NOT FOUND *****		
30	1,2-Dichloroethane	***** NOT FOUND *****		
31	1,1-Dichloropropene	***** NOT FOUND *****		
32	Benzene	***** NOT FOUND *****		

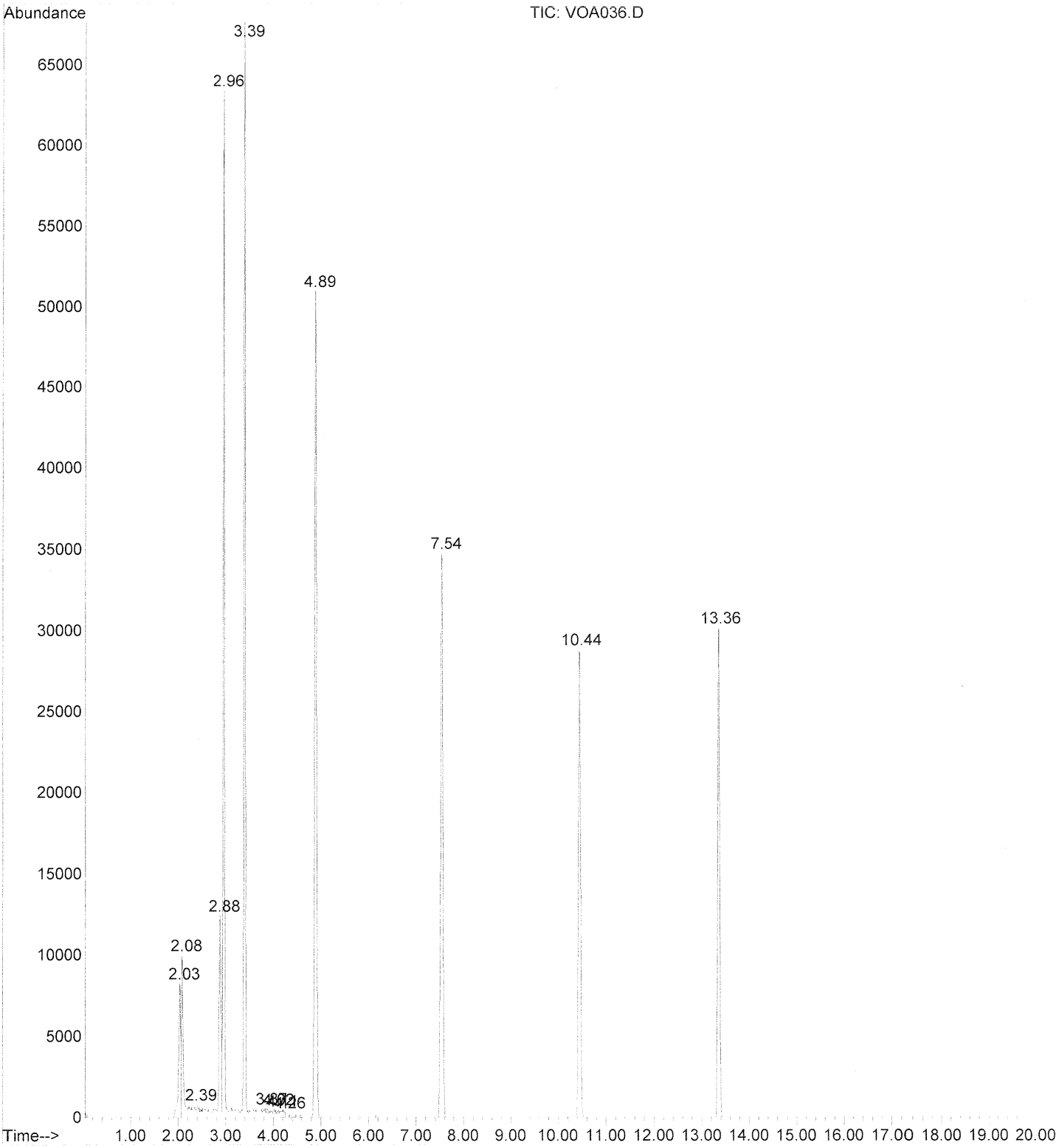
33	Carbon Tetrachloride	***** NOT FOUND *****		
34	2-Chloroethanol	***** NOT FOUND *****		
35	1,4-Difluorobenzene	3.39	100.0 ppb	ISTD2#
36	1,2-Dichloropropane	***** NOT FOUND *****		
37	Trichloroethene	***** NOT FOUND *****		
38	Dibromomethane	***** NOT FOUND *****		
39	Bromodichloromethane	***** NOT FOUND *****		
40	Methyl methacrylate	***** NOT FOUND *****		
41	1,4-Dioxane	***** NOT FOUND *****		
42	cis-1,3-Dichloropropene	***** NOT FOUND *****		
43	2-Nitropropane	***** NOT FOUND *****		
44	trans-1,3-Dichloropropene	***** NOT FOUND *****		
45	Toluene	***** NOT FOUND *****		
46	Toluene-d8-surrogate	4.90	84.5 ppb	85 LR
47	1,1,2-Trichloroethane	***** NOT FOUND *****		
48	1,3-Dichloropropane	***** NOT FOUND *****		
49	Ethylmethacrylate	***** NOT FOUND *****		
50	Dibromochloromethane	***** NOT FOUND *****		
51	1,2-Dibromomethane	***** NOT FOUND *****		
52	Tetrachloroethene	***** NOT FOUND *****		
53	Chlorobenzene	***** NOT FOUND *****		
54	Chlorobenzene-d5	7.55	100.0 ppb	ISTD3
55	1,1,1,2-Tetrachloroethane	***** NOT FOUND *****		
56	Ethylbenzene	***** NOT FOUND *****		
57	m/p-xylene	***** NOT FOUND *****		
58	Bromoform	***** NOT FOUND *****		
59	Styrene	***** NOT FOUND *****		
60	o-Xylene	***** NOT FOUND *****		
61	cis-1,4-Dichloro-2-butene	***** NOT FOUND *****		
62	trans-1,4-Dichloro-2-butene	***** NOT FOUND *****		
63	Bromofluorobenzene-surrogate	10.44	101.7 ppb	102 LR
64	1,1,2,2-Tetrachloroethane	***** NOT FOUND *****		
65	1,2,3-Trichloropropane	***** NOT FOUND *****		
66	Isopropylbenzene	***** NOT FOUND *****		
67	Bromobenzene	***** NOT FOUND *****		
68	2-Chlorotoluene	***** NOT FOUND *****		
69	n-Propylbenzene	***** NOT FOUND *****		
70	4-Chlorotoluene	***** NOT FOUND *****		
71	1,3,5-Trimethylbenzene	***** NOT FOUND *****		
72	Pentachloroethane	***** NOT FOUND *****		
73	tert-Butylbenzene	***** NOT FOUND *****		
74	1,2,4-Trimethylbenzene	***** NOT FOUND *****		
75	1,3-Dichlorobenzene	***** NOT FOUND *****		
76	1,4-Dichlorobenzene-d4	13.36	100.0 ppb	ISTD4
77	sec-Butylbenzene	***** NOT FOUND *****		
78	1,4-Dichlorobenzene	***** NOT FOUND *****		
79	p-iso-Propyltoluene	***** NOT FOUND *****		
80	1,2-Dichlorobenzene	***** NOT FOUND *****		
81	n-Butylbenzene	***** NOT FOUND *****		
82	1,2-Bromo-3-chloropropane	***** NOT FOUND *****		
83	Nitobenzene	***** NOT FOUND *****		
84	1,2,4-Trichlorobenzene	***** NOT FOUND *****		
85	Naphthalene	***** NOT FOUND *****		
86	1,2,3-Trichlorobenzene	***** NOT FOUND *****		
87	Hexachlorobutadiene	***** NOT FOUND *****		

END OF REPORT

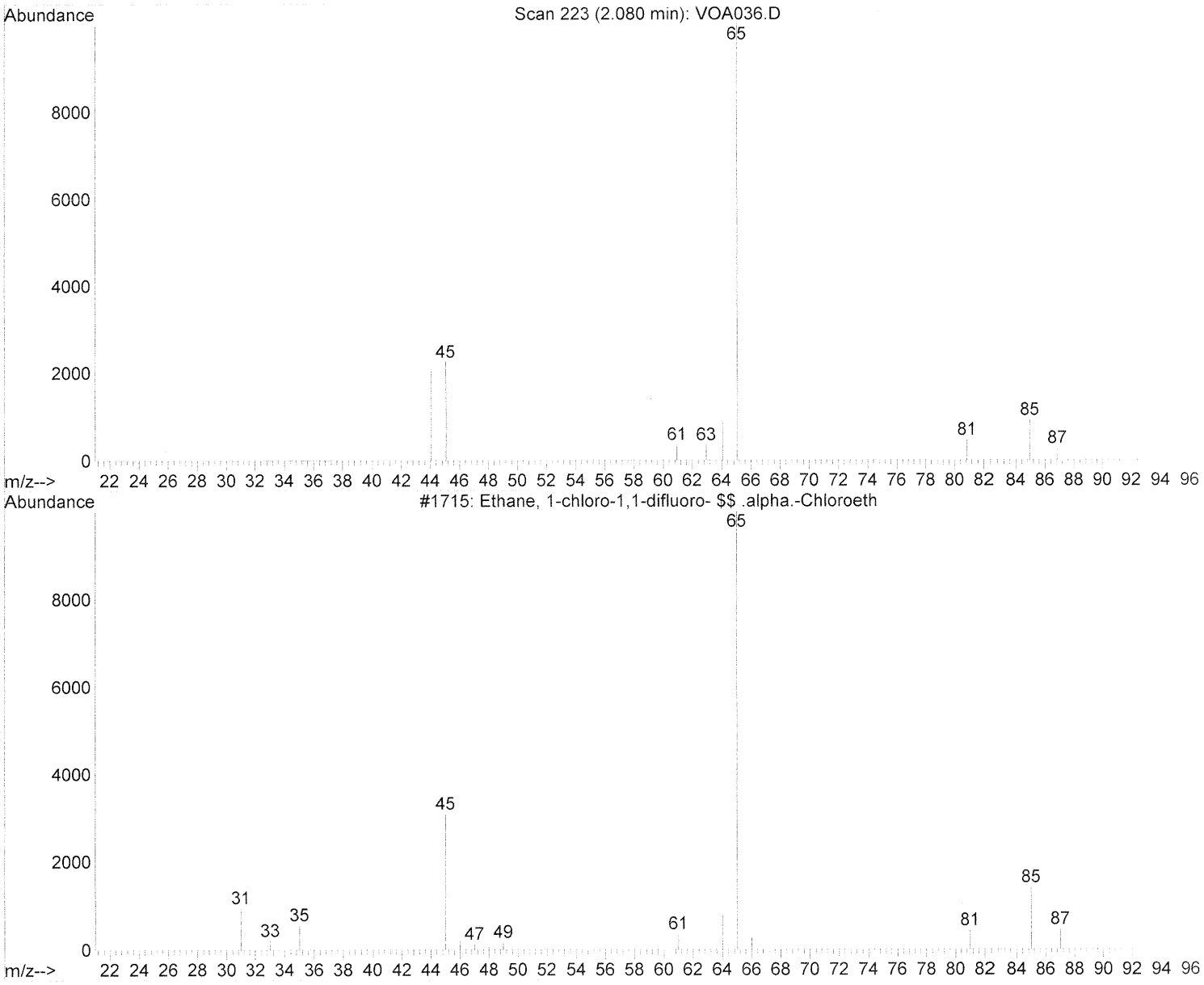
Qualifiers Not Satisfied

Fri May 26 12:29:47 2006

File : C:\HPCHEM\1\DATA\052406\VOA036.D
Operator : dmt
Acquired : 26 May 2006 12:57 pm using AcqMethod VOA0506
Instrument : GCD Syste
Sample Name: MW-3 H2O
Misc Info :
Vial Number: 1



Library Searched : C:\DATABASE\NIST129K.L
Quality : 83
ID : Ethane, 1-chloro-1,1-difluoro- $$ .alpha.-Chloroethylidene fluoride $$ Freon 142 $$ Freon 142b $$ FC 142b $$ Genetron 101 $$ Genetro



Calibration Last Updated: Fri May 26 08:11:46 2006
Quantitation Report - Summary

Information from Data File:
File : C:\HPCHEM\1\DATA\052406\VOA036.D
Operator : dmt
Acquired : 26 May 2006 12:57 pm using AcqMethod VOA0506
Sample Name: MW-3 H2O
Misc Info :
Vial Number: 1
CurrentMeth: C:\HPCHEM\1\METHODS\VOA0506R.M

Quantitation Settings:
Non-Reference Peak Window: 10.00 Percent
Correlation Window: 0.02 Minutes
Default Multiplier: 1.00
Default Sample Concentration: 0.00
Peak Type Decoding: m -> Manually re-integrated

Number of Compounds in Database: 87

#	Compound	Ret Time	Concentration	Peak Type
1	Pentafluorobenzene	2.96	100.0 ppb	ISTD1
2	Dichlorofluoromethane	2.08	3.0 ppb	#
3	Chloromethane	***** NOT FOUND *****		
4	Vinyl Chloride	***** NOT FOUND *****		
5	Bromomethane	***** NOT FOUND *****		
6	Chloroethane	2.08	6.0 ppb	#
7	Trichlorofluoromethane	***** NOT FOUND *****		
8	Acetonitrile	***** NOT FOUND *****		
9	Diethyl ether	***** NOT FOUND *****		
10	1,1-Dichloroethene	***** NOT FOUND *****		
11	Acrylonitrile	***** NOT FOUND *****		
12	1,1,2-Trichlorotrifluoroethane	***** NOT FOUND *****		
13	Methylene Chloride	2.38	4.0 ppb	#
14	Iodomethane	***** NOT FOUND *****		
15	trans-1,2-Dichloroethene	***** NOT FOUND *****		
16	MTBE	***** NOT FOUND *****		
17	Allyl chloride	***** NOT FOUND *****		
18	1,1-Dichloroethane	***** NOT FOUND *****		
19	Carbon disulfide	***** NOT FOUND *****		
20	Chloroprene	***** NOT FOUND *****		
21	Methacrylonitrile	***** NOT FOUND *****		
22	cis-1,2-Dichloroethene	***** NOT FOUND *****		
23	Methyl acrylate	***** NOT FOUND *****		
24	Chloroform	***** NOT FOUND *****		
25	2,2-Dichloropropane	***** NOT FOUND *****		
26	Dibromofluoromethane-surrogate	2.88	82.0 ppb	
27	Isobutyl alcohol	***** NOT FOUND *****		
28	Tetrahydrofuran	***** NOT FOUND *****		
29	1,1,1-Trichloroethane	***** NOT FOUND *****		
30	1,2-Dichloroethane	***** NOT FOUND *****		
31	1,1-Dichloropropene	***** NOT FOUND *****		
32	Benzene	***** NOT FOUND *****		

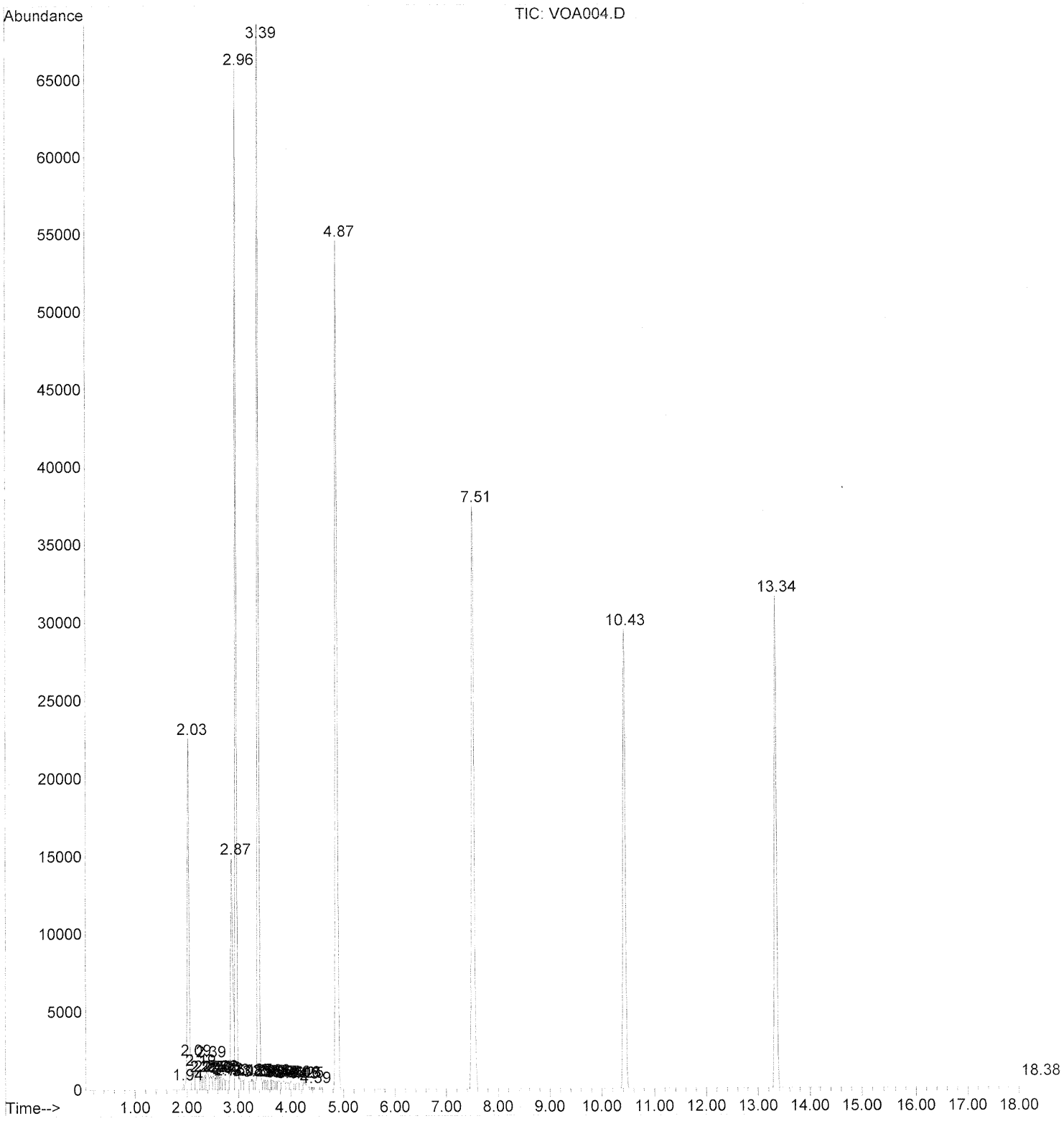
33 Carbon Tetrachloride	***** NOT FOUND *****		
34 2-Chloroethanol	***** NOT FOUND *****		
35 1,4-Difluorobenzene	3.39	100.0 ppb	ISTD2
36 1,2-Dichloropropane	***** NOT FOUND *****		
37 Trichloroethene	***** NOT FOUND *****		
38 Dibromomethane	***** NOT FOUND *****		
39 Bromodichloromethane	***** NOT FOUND *****		
40 Methyl methacrylate	***** NOT FOUND *****		
41 1,4-Dioxane	***** NOT FOUND *****		
42 cis-1,3-Dichloropropene	***** NOT FOUND *****		
43 2-Nitropropane	***** NOT FOUND *****		
44 trans-1,3-Dichloropropene	***** NOT FOUND *****		
45 Toluene	***** NOT FOUND *****		
46 Toluene-d8-surrogate	4.89	97.5 ppb	98 %
47 1,1,2-Trichloroethane	***** NOT FOUND *****		
48 1,3-Dichloropropane	***** NOT FOUND *****		
49 Ethylmethacrylate	***** NOT FOUND *****		
50 Dibromochloromethane	***** NOT FOUND *****		
51 1,2-Dibromomethane	***** NOT FOUND *****		
52 Tetrachloroethene	***** NOT FOUND *****		
53 Chlorobenzene	***** NOT FOUND *****		
54 Chlorobenzene-d5	7.54	100.0 ppb	ISTD3
55 1,1,1,2-Tetrachloroethane	***** NOT FOUND *****		
56 Ethylbenzene	***** NOT FOUND *****		
57 m/p-xylene	***** NOT FOUND *****		
58 Bromoform	***** NOT FOUND *****		
59 Styrene	***** NOT FOUND *****		
60 o-Xylene	***** NOT FOUND *****		
61 cis-1,4-Dichloro-2-butene	***** NOT FOUND *****		
62 trans-1,4-Dichloro-2-butene	***** NOT FOUND *****		
63 Bromofluorobenzene-surrogate	10.44	94.8 ppb	95 %
64 1,1,2,2-Tetrachloroethane	***** NOT FOUND *****		
65 1,2,3-Trichloropropane	***** NOT FOUND *****		
66 Isopropylbenzene	***** NOT FOUND *****		
67 Bromobenzene	***** NOT FOUND *****		
68 2-Chlorotoluene	***** NOT FOUND *****		
69 n-Propylbenzene	***** NOT FOUND *****		
70 4-Chlorotoluene	***** NOT FOUND *****		
71 1,3,5-Trimethylbenzene	***** NOT FOUND *****		
72 Pentachloroethane	***** NOT FOUND *****		
73 tert-Butylbenzene	***** NOT FOUND *****		
74 1,2,4-Trimethylbenzene	***** NOT FOUND *****		
75 1,3-Dichlorobenzene	***** NOT FOUND *****		
76 1,4-Dichlorobenzene-d4	13.36	100.0 ppb	ISTD4
77 sec-Butylbenzene	***** NOT FOUND *****		
78 1,4-Dichlorobenzene	***** NOT FOUND *****		
79 p-iso-Propyltoluene	***** NOT FOUND *****		
80 1,2-Dichlorobenzene	***** NOT FOUND *****		
81 n-Butylbenzene	***** NOT FOUND *****		
82 1,2-Bromo-3-chloropropane	***** NOT FOUND *****		
83 Nitobenzene	***** NOT FOUND *****		
84 1,2,4-Trichlorobenzene	***** NOT FOUND *****		
85 Naphthalene	***** NOT FOUND *****		
86 1,2,3-Trichlorobenzene	***** NOT FOUND *****		
87 Hexachlorobutadiene	***** NOT FOUND *****		

END OF REPORT

Qualifiers Not Satisfied

Fri May 26 13:23:32 2006

File : C:\HPCHEM\1\DATA\052406\VOA004.D
Operator : dmt
Acquired : 24 May 2006 11:26 am using AcqMethod VOA0506
Instrument : GCD Syste
Sample Name: MTH BLK 5/24
Misc Info :
Vial Number: 1



Calibration Last Updated: Wed May 24 11:22:20 2006
Quantitation Report - Summary

Information from Data File:
File : C:\HPCHEM\1\DATA\052406\VOA004.D
Operator : dmt
Acquired : 24 May 2006 11:26 am using AcqMethod VOA0506
Sample Name: MTH BLK 5/24
Misc Info :
Vial Number: 1
CurrentMeth: C:\HPCHEM\1\METHODS\VOA0506R.M

Quantitation Settings:
Non-Reference Peak Window: 10.00 Percent
Correlation Window: 0.02 Minutes
Default Multiplier: 1.00
Default Sample Concentration: 0.00
Peak Type Decoding: m -> Manually re-integrated

Number of Compounds in Database: 87

#	Compound	Ret Time	Concentration	Peak Type
1	Pentafluorobenzene	2.96	100.0 ppb	ISTD1
2	Dichlorofluoromethane	***** NOT FOUND *****		
3	Chloromethane	***** NOT FOUND *****		
4	Vinyl Chloride	***** NOT FOUND *****		
5	Bromomethane	***** NOT FOUND *****		
6	Chloroethane	***** NOT FOUND *****		
7	Trichlorofluoromethane	***** NOT FOUND *****		
8	Acetonitrile	***** NOT FOUND *****		
9	Diethyl ether	***** NOT FOUND *****		
10	1,1-Dichloroethene	***** NOT FOUND *****		
11	Acrylonitrile	***** NOT FOUND *****		
12	1,1,2-Trichlorotrifluoroethane	***** NOT FOUND *****		
13	Methylene Chloride	2.39	5.7 ppb	<
14	Iodomethane	***** NOT FOUND *****		
15	trans-1,2-Dichloroethene	***** NOT FOUND *****		
16	MTBE	***** NOT FOUND *****		
17	Allyl chloride	***** NOT FOUND *****		
18	1,1-Dichloroethane	***** NOT FOUND *****		
19	Carbon disulfide	***** NOT FOUND *****		
20	Chloroprene	***** NOT FOUND *****		
21	Methacrylonitrile	***** NOT FOUND *****		
22	cis-1,2-Dichloroethene	***** NOT FOUND *****		
23	Methyl acrylate	***** NOT FOUND *****		
24	Chloroform	***** NOT FOUND *****		
25	2,2-Dichloropropane	***** NOT FOUND *****		
26	Dibromofluoromethane-surrogate	2.87	100.8 ppb	101 %R
27	Isobutyl alcohol	***** NOT FOUND *****		
28	Tetrahydrofuran	***** NOT FOUND *****		
29	1,1,1-Trichloroethane	***** NOT FOUND *****		
30	1,2-Dichloroethane	***** NOT FOUND *****		
31	1,1-Dichloropropene	***** NOT FOUND *****		
32	Benzene	***** NOT FOUND *****		

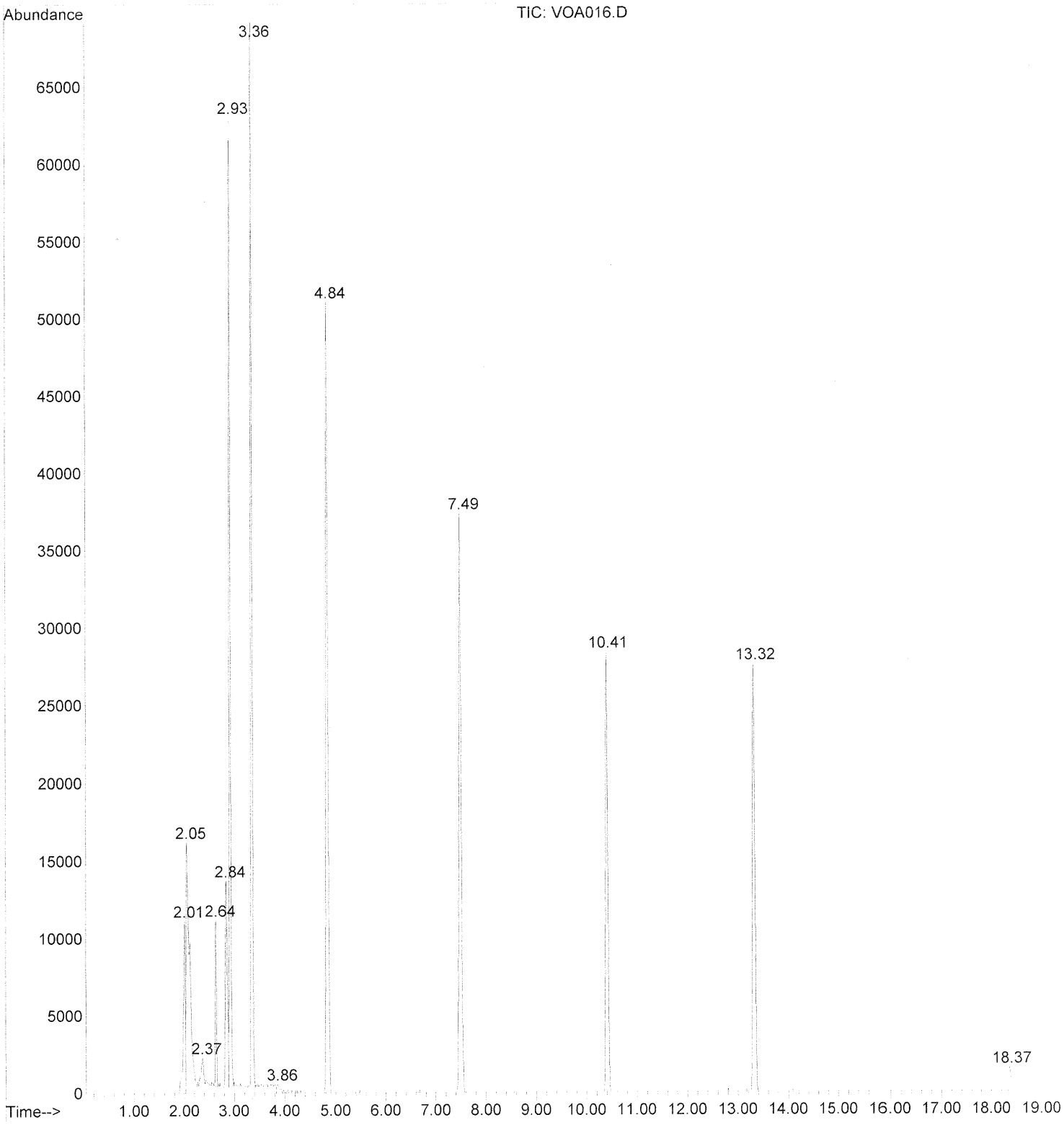
33	Carbon Tetrachloride	***** NOT FOUND *****		
34	2-Chloroethanol	***** NOT FOUND *****		
35	1,4-Difluorobenzene	3.39	100.0 ppb	ISTD2
36	1,2-Dichloropropane	***** NOT FOUND *****		
37	Trichloroethene	***** NOT FOUND *****		
38	Dibromomethane	***** NOT FOUND *****		
39	Bromodichloromethane	***** NOT FOUND *****		
40	Methyl methacrylate	***** NOT FOUND *****		
41	1,4-Dioxane	***** NOT FOUND *****		
42	cis-1,3-Dichloropropene	***** NOT FOUND *****		
43	2-Nitropropane	***** NOT FOUND *****		
44	trans-1,3-Dichloropropene	***** NOT FOUND *****		
45	Toluene	***** NOT FOUND *****		
46	Toluene-d8-surrogate	4.87	96.2 ppb	96 %R
47	1,1,2-Trichloroethane	***** NOT FOUND *****		
48	1,3-Dichloropropane	***** NOT FOUND *****		
49	Ethylmethacrylate	***** NOT FOUND *****		
50	Dibromochloromethane	***** NOT FOUND *****		
51	1,2-Dibromomethane	***** NOT FOUND *****		
52	Tetrachloroethene	***** NOT FOUND *****		
53	Chlorobenzene	***** NOT FOUND *****		
54	Chlorobenzene-d5	7.51	100.0 ppb	ISTD3
55	1,1,1,2-Tetrachloroethane	***** NOT FOUND *****		
56	Ethylbenzene	***** NOT FOUND *****		
57	m/p-xylene	***** NOT FOUND *****		
58	Bromoform	***** NOT FOUND *****		
59	Styrene	***** NOT FOUND *****		
60	o-Xylene	***** NOT FOUND *****		
61	cis-1,4-Dichloro-2-butene	***** NOT FOUND *****		
62	trans-1,4-Dichloro-2-butene	***** NOT FOUND *****		
63	Bromofluorobenzene-surrogate	10.43	95.1 ppb	95 %R
64	1,1,2,2-Tetrachloroethane	***** NOT FOUND *****		
65	1,2,3-Trichloropropane	***** NOT FOUND *****		
66	Isopropylbenzene	***** NOT FOUND *****		
67	Bromobenzene	***** NOT FOUND *****		
68	2-Chlorotoluene	***** NOT FOUND *****		
69	n-Propylbenzene	***** NOT FOUND *****		
70	4-Chlorotoluene	***** NOT FOUND *****		
71	1,3,5-Trimethylbenzene	***** NOT FOUND *****		
72	Pentachloroethane	***** NOT FOUND *****		
73	tert-Butylbenzene	***** NOT FOUND *****		
74	1,2,4-Trimethylbenzene	***** NOT FOUND *****		
75	1,3-Dichlorobenzene	***** NOT FOUND *****		
76	1,4-Dichlorobenzene-d4	13.34	100.0 ppb	ISTD4
77	sec-Butylbenzene	***** NOT FOUND *****		
78	1,4-Dichlorobenzene	***** NOT FOUND *****		
79	p-iso-Propyltoluene	***** NOT FOUND *****		
80	1,2-Dichlorobenzene	***** NOT FOUND *****		
81	n-Butylbenzene	***** NOT FOUND *****		
82	1,2-Bromo-3-chloropropane	***** NOT FOUND *****		
83	Nitobenzene	***** NOT FOUND *****		
84	1,2,4-Trichlorobenzene	***** NOT FOUND *****		
85	Naphthalene	18.38	22.7 ppb	#
86	1,2,3-Trichlorobenzene	***** NOT FOUND *****		
87	Hexachlorobutadiene	***** NOT FOUND *****		

END OF REPORT

Qualifiers Not Satisfied

Wed May 24 11:45:12 2006

File : C:\HPCHEM\1\DATA\052406\VOA016.D
Operator : dmt
Acquired : 25 May 2006 8:31 am using AcqMethod VOA0506
Instrument : GCD Syste
Sample Name: MTH BLK 5/25 SOIL
Misc Info :
Vial Number: 1



Calibration Last Updated: Thu May 25 08:53:45 2006
Quantitation Report - Summary

Information from Data File:
File : C:\HPCHEM\1\DATA\052406\VOA016.D
Operator : dmt
Acquired : 25 May 2006 8:31 am using AcqMethod VOA0506
Sample Name: MTH BLK 5/25 SOIL
Misc Info :
Vial Number: 1
CurrentMeth: C:\HPCHEM\1\METHODS\VOA0506R.M

Quantitation Settings:
Non-Reference Peak Window: 10.00 Percent
Correlation Window: 0.02 Minutes
Default Multiplier: 1.00
Default Sample Concentration: 0.00
Peak Type Decoding: m -> Manually re-integrated

Number of Compounds in Database: 87

#	Compound	Ret Time	Concentration	Peak Type
1	Pentafluorobenzene	2.93	100.0 ppb	ISTD1
2	Dichlorofluoromethane	***** NOT FOUND *****		
3	Chloromethane	***** NOT FOUND *****		
4	Vinyl Chloride	***** NOT FOUND *****		
5	Bromomethane	***** NOT FOUND *****		
6	Chloroethane	***** NOT FOUND *****		
7	Trichlorofluoromethane	***** NOT FOUND *****		
8	Acetonitrile	***** NOT FOUND *****		
9	Diethyl ether	***** NOT FOUND *****		
10	1,1-Dichloroethene	***** NOT FOUND *****		
11	Acrylonitrile	***** NOT FOUND *****		
12	1,1,2-Trichlorotrifluoroethane	***** NOT FOUND *****		
13	Methylene Chloride	***** NOT FOUND *****		
14	Iodomethane	***** NOT FOUND *****		
15	trans-1,2-Dichloroethene	***** NOT FOUND *****		
16	MTBE	***** NOT FOUND *****		
17	Allyl chloride	***** NOT FOUND *****		
18	1,1-Dichloroethane	***** NOT FOUND *****		
19	Carbon disulfide	***** NOT FOUND *****		
20	Chloroprene	***** NOT FOUND *****		
21	Methacrylonitrile	***** NOT FOUND *****		
22	cis-1,2-Dichloroethene	***** NOT FOUND *****		
23	Methyl acrylate	***** NOT FOUND *****		
24	Chloroform	***** NOT FOUND *****		
25	2,2-Dichloropropane	***** NOT FOUND *****		
26	Dibromofluoromethane-surrogate	2.83	98.1 ppb	
27	Isobutyl alcohol	***** NOT FOUND *****		
28	Tetrahydrofuran	***** NOT FOUND *****		
29	1,1,1-Trichloroethane	***** NOT FOUND *****		
30	1,2-Dichloroethane	***** NOT FOUND *****		
31	1,1-Dichloropropene	***** NOT FOUND *****		
32	Benzene	***** NOT FOUND *****		

98 1/R

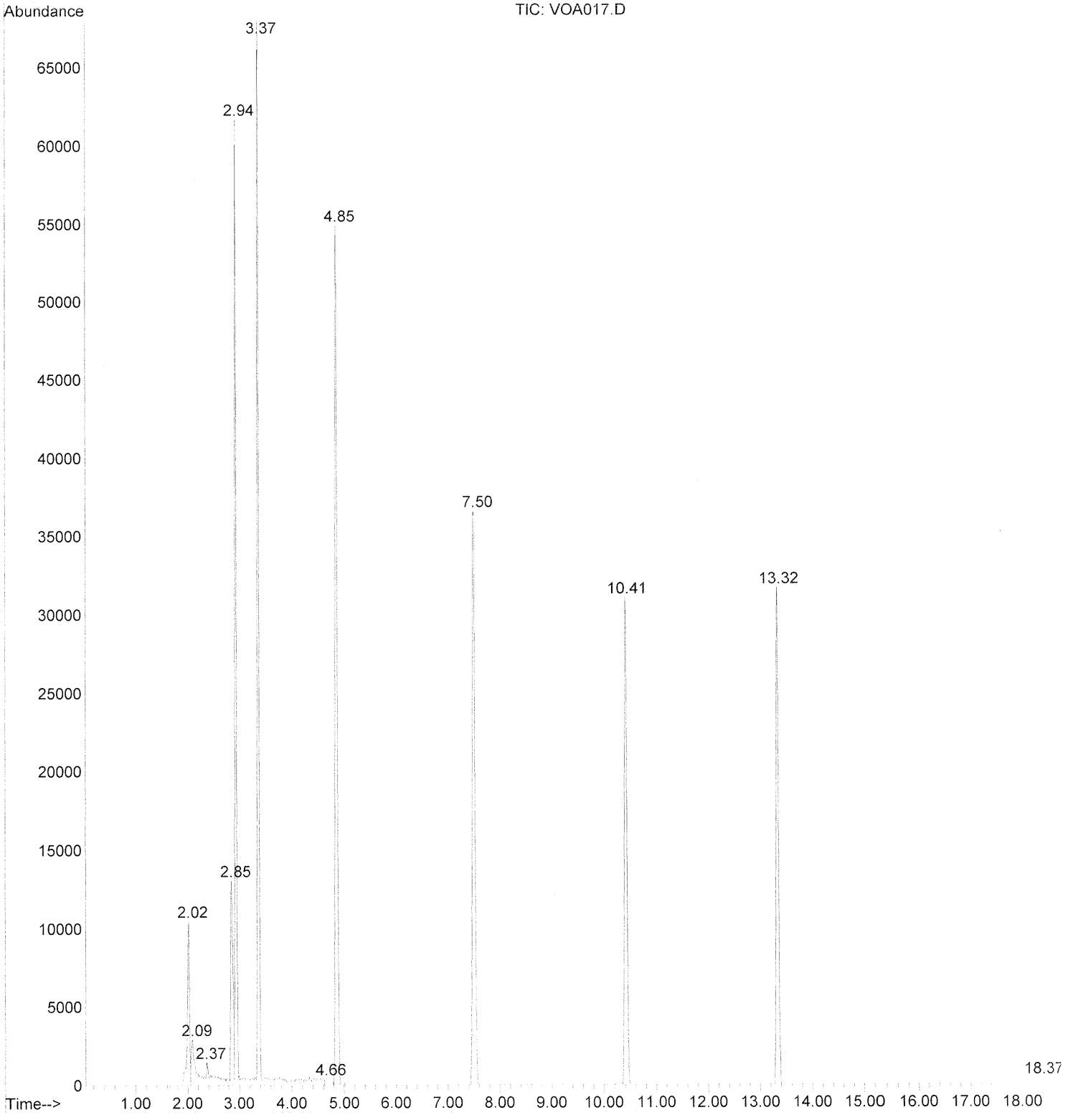
33 Carbon Tetrachloride	***** NOT FOUND *****		
34 2-Chloroethanol	***** NOT FOUND *****		
35 1,4-Difluorobenzene	3.36	100.0 ppb	ISTD2
36 1,2-Dichloropropane	***** NOT FOUND *****		
37 Trichloroethene	***** NOT FOUND *****		
38 Dibromomethane	***** NOT FOUND *****		
39 Bromodichloromethane	***** NOT FOUND *****		
40 Methyl methacrylate	***** NOT FOUND *****		
41 1,4-Dioxane	***** NOT FOUND *****		
42 cis-1,3-Dichloropropene	***** NOT FOUND *****		
43 2-Nitropropane	***** NOT FOUND *****		
44 trans-1,3-Dichloropropene	***** NOT FOUND *****		
45 Toluene	***** NOT FOUND *****		
46 Toluene-d8-surrogate	4.84	89.7 ppb	
47 1,1,2-Trichloroethane	***** NOT FOUND *****		
48 1,3-Dichloropropane	***** NOT FOUND *****		
49 Ethylmethacrylate	***** NOT FOUND *****		
50 Dibromochloromethane	***** NOT FOUND *****		
51 1,2-Dibromomethane	***** NOT FOUND *****		
52 Tetrachloroethene	***** NOT FOUND *****		
53 Chlorobenzene	***** NOT FOUND *****		
54 Chlorobenzene-d5	7.49	100.0 ppb	ISTD3
55 1,1,1,2-Tetrachloroethane	***** NOT FOUND *****		
56 Ethylbenzene	***** NOT FOUND *****		
57 m/p-xylene	***** NOT FOUND *****		
58 Bromoform	***** NOT FOUND *****		
59 Styrene	***** NOT FOUND *****		
60 o-Xylene	***** NOT FOUND *****		
61 cis-1,4-Dichloro-2-butene	***** NOT FOUND *****		
62 trans-1,4-Dichloro-2-butene	***** NOT FOUND *****		
63 Bromofluorobenzene-surrogate	10.41	89.8 ppb	
64 1,1,2,2-Tetrachloroethane	***** NOT FOUND *****		
65 1,2,3-Trichloropropane	***** NOT FOUND *****		
66 Isopropylbenzene	***** NOT FOUND *****		
67 Bromobenzene	***** NOT FOUND *****		
68 2-Chlorotoluene	***** NOT FOUND *****		
69 n-Propylbenzene	***** NOT FOUND *****		
70 4-Chlorotoluene	***** NOT FOUND *****		
71 1,3,5-Trimethylbenzene	***** NOT FOUND *****		
72 Pentachloroethane	***** NOT FOUND *****		
73 tert-Butylbenzene	***** NOT FOUND *****		
74 1,2,4-Trimethylbenzene	***** NOT FOUND *****		
75 1,3-Dichlorobenzene	***** NOT FOUND *****		
76 1,4-Dichlorobenzene-d4	13.33	100.0 ppb	ISTD4
77 sec-Butylbenzene	***** NOT FOUND *****		
78 1,4-Dichlorobenzene	***** NOT FOUND *****		
79 p-iso-Propyltoluene	***** NOT FOUND *****		
80 1,2-Dichlorobenzene	***** NOT FOUND *****		
81 n-Butylbenzene	***** NOT FOUND *****		
82 1,2-Bromo-3-chloropropane	***** NOT FOUND *****		
83 Nitobenzene	***** NOT FOUND *****		
84 1,2,4-Trichlorobenzene	***** NOT FOUND *****		
85 Naphthalene	18.37	53.0 ppb	#
86 1,2,3-Trichlorobenzene	***** NOT FOUND *****		
87 Hexachlorobutadiene	***** NOT FOUND *****		

END OF REPORT

Qualifiers Not Satisfied

Thu May 25 08:53:52 2006

File : C:\HPCHEM\1\DATA\052406\VOA017.D
Operator : dmt
Acquired : 25 May 2006 8:57 am using AcqMethod VOA0506
Instrument : GCD Syste
Sample Name: MTH BLK 5/25 H2O
Misc Info :
Vial Number: 1



Calibration Last Updated: Thu May 25 08:53:45 2006
Quantitation Report - Summary

Information from Data File:
File : C:\HPCHEM\1\DATA\052406\VOA017.D
Operator : dmt
Acquired : 25 May 2006 8:57 am using AcqMethod VOA0506
Sample Name: MTH BLK 5/25 H2O
Misc Info :
Vial Number: 1
CurrentMeth: C:\HPCHEM\1\METHODS\VOA0506R.M

Quantitation Settings:
Non-Reference Peak Window: 10.00 Percent
Correlation Window: 0.02 Minutes
Default Multiplier: 1.00
Default Sample Concentration: 0.00
Peak Type Decoding: m -> Manually re-integrated

Number of Compounds in Database: 87

#	Compound	Ret Time	Concentration	Peak Type
1	Pentafluorobenzene	2.94	100.0 ppb	ISTD1
2	Dichlorofluoromethane	***** NOT FOUND *****		
3	Chloromethane	***** NOT FOUND *****		
4	Vinyl Chloride	***** NOT FOUND *****		
5	Bromomethane	***** NOT FOUND *****		
6	Chloroethane	***** NOT FOUND *****		
7	Trichlorofluoromethane	***** NOT FOUND *****		
8	Acetonitrile	***** NOT FOUND *****		
9	Diethyl ether	***** NOT FOUND *****		
10	1,1-Dichloroethene	***** NOT FOUND *****		
11	Acrylonitrile	***** NOT FOUND *****		
12	1,1,2-Trichlorotrifluoroethane	***** NOT FOUND *****		
13	Methylene Chloride	2.37	7.5 ppb	#
14	Iodomethane	***** NOT FOUND *****		
15	trans-1,2-Dichloroethene	***** NOT FOUND *****		
16	MTBE	***** NOT FOUND *****		
17	Allyl chloride	***** NOT FOUND *****		
18	1,1-Dichloroethane	***** NOT FOUND *****		
19	Carbon disulfide	***** NOT FOUND *****		
20	Chloroprene	***** NOT FOUND *****		
21	Methacrylonitrile	***** NOT FOUND *****		
22	cis-1,2-Dichloroethene	***** NOT FOUND *****		
23	Methyl acrylate	***** NOT FOUND *****		
24	Chloroform	***** NOT FOUND *****		
25	2,2-Dichloropropane	***** NOT FOUND *****		
26	Dibromofluoromethane-surrogate	2.84	93.1 ppb	
27	Isobutyl alcohol	***** NOT FOUND *****		
28	Tetrahydrofuran	***** NOT FOUND *****		
29	1,1,1-Trichloroethane	***** NOT FOUND *****		
30	1,2-Dichloroethane	***** NOT FOUND *****		
31	1,1-Dichloropropene	***** NOT FOUND *****		
32	Benzene	***** NOT FOUND *****		

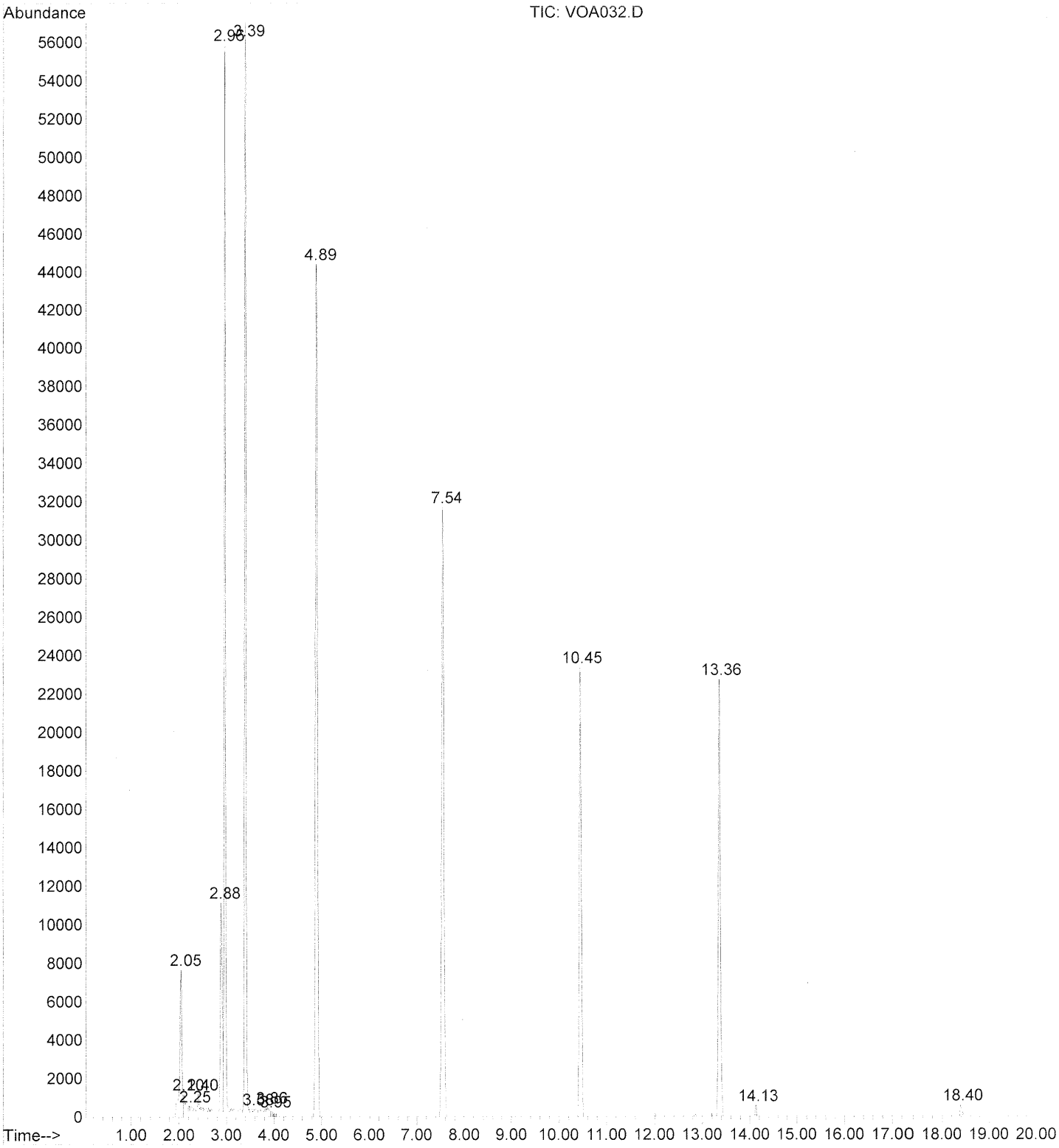
33 Carbon Tetrachloride	***** NOT FOUND *****		
34 2-Chloroethanol	***** NOT FOUND *****		
35 1,4-Difluorobenzene	3.37	100.0 ppb	ISTD2
36 1,2-Dichloropropane	***** NOT FOUND *****		
37 Trichloroethene	***** NOT FOUND *****		
38 Dibromomethane	***** NOT FOUND *****		
39 Bromodichloromethane	***** NOT FOUND *****		
40 Methyl methacrylate	***** NOT FOUND *****		
41 1,4-Dioxane	***** NOT FOUND *****		
42 cis-1,3-Dichloropropene	***** NOT FOUND *****		
43 2-Nitropropane	***** NOT FOUND *****		
44 trans-1,3-Dichloropropene	***** NOT FOUND *****		
45 Toluene	***** NOT FOUND *****		
46 Toluene-d8-surrogate	4.85	109.4 ppb	109 %R
47 1,1,2-Trichloroethane	***** NOT FOUND *****		
48 1,3-Dichloropropane	***** NOT FOUND *****		
49 Ethylmethacrylate	***** NOT FOUND *****		
50 Dibromochloromethane	***** NOT FOUND *****		
51 1,2-Dibromomethane	***** NOT FOUND *****		
52 Tetrachloroethene	***** NOT FOUND *****		
53 Chlorobenzene	***** NOT FOUND *****		
54 Chlorobenzene-d5	7.50	100.0 ppb	ISTD3
55 1,1,1,2-Tetrachloroethane	***** NOT FOUND *****		
56 Ethylbenzene	***** NOT FOUND *****		
57 m/p-xylene	***** NOT FOUND *****		
58 Bromoform	***** NOT FOUND *****		
59 Styrene	***** NOT FOUND *****		
60 o-Xylene	***** NOT FOUND *****		
61 cis-1,4-Dichloro-2-butene	***** NOT FOUND *****		
62 trans-1,4-Dichloro-2-butene	***** NOT FOUND *****		
63 Bromofluorobenzene-surrogate	10.42	103.8 ppb	104 %R
64 1,1,2,2-Tetrachloroethane	***** NOT FOUND *****		
65 1,2,3-Trichloropropane	***** NOT FOUND *****		
66 Isopropylbenzene	***** NOT FOUND *****		
67 Bromobenzene	***** NOT FOUND *****		
68 2-Chlorotoluene	***** NOT FOUND *****		
69 n-Propylbenzene	***** NOT FOUND *****		
70 4-Chlorotoluene	***** NOT FOUND *****		
71 1,3,5-Trimethylbenzene	***** NOT FOUND *****		
72 Pentachloroethane	***** NOT FOUND *****		
73 tert-Butylbenzene	***** NOT FOUND *****		
74 1,2,4-Trimethylbenzene	***** NOT FOUND *****		
75 1,3-Dichlorobenzene	***** NOT FOUND *****		
76 1,4-Dichlorobenzene-d4	13.33	100.0 ppb	ISTD4
77 sec-Butylbenzene	***** NOT FOUND *****		
78 1,4-Dichlorobenzene	***** NOT FOUND *****		
79 p-iso-Propyltoluene	***** NOT FOUND *****		
80 1,2-Dichlorobenzene	***** NOT FOUND *****		
81 n-Butylbenzene	***** NOT FOUND *****		
82 1,2-Bromo-3-chloropropane	***** NOT FOUND *****		
83 Nitobenzene	***** NOT FOUND *****		
84 1,2,4-Trichlorobenzene	***** NOT FOUND *****		
85 Naphthalene	18.37	19.8 ppb	#
86 1,2,3-Trichlorobenzene	***** NOT FOUND *****		
87 Hexachlorobutadiene	***** NOT FOUND *****		

END OF REPORT

Qualifiers Not Satisfied

Thu May 25 09:16:23 2006

File : C:\HPCHEM\1\DATA\052406\VOA032.D
Operator : dmt
Acquired : 26 May 2006 8:18 am using AcqMethod VOA0506
Instrument : GCD Syste
Sample Name: MTH BLK 5/26 H2O
Misc Info :
Vial Number: 1



Calibration Last Updated: Fri May 26 08:11:46 2006
Quantitation Report - Summary

Information from Data File:
File : C:\HPCHEM\1\DATA\052406\VOA032.D
Operator : dmt
Acquired : 26 May 2006 8:18 am using AcqMethod VOA0506
Sample Name: MTH BLK 5/26 H2O
Misc Info :
Vial Number: 1
CurrentMeth: C:\HPCHEM\1\METHODS\VOA0506R.M

Quantitation Settings:
Non-Reference Peak Window: 10.00 Percent
Correlation Window: 0.02 Minutes
Default Multiplier: 1.00
Default Sample Concentration: 0.00
Peak Type Decoding: m -> Manually re-integrated

Number of Compounds in Database: 87

#	Compound	Ret Time	Concentration	Peak Type
1	Pentafluorobenzene	2.96	100.0 ppb	ISTD1
2	Dichlorofluoromethane	***** NOT FOUND *****		
3	Chloromethane	***** NOT FOUND *****		
4	Vinyl Chloride	***** NOT FOUND *****		
5	Bromomethane	***** NOT FOUND *****		
6	Chloroethane	***** NOT FOUND *****		
7	Trichlorofluoromethane	***** NOT FOUND *****		
8	Acetonitrile	***** NOT FOUND *****		
9	Diethyl ether	***** NOT FOUND *****		
10	1,1-Dichloroethene	***** NOT FOUND *****		
11	Acrylonitrile	***** NOT FOUND *****		
12	1,1,2-Trichlorotrifluoroethane	***** NOT FOUND *****		
13	Methylene Chloride	2.40	7.1 ppb	#
14	Iodomethane	***** NOT FOUND *****		
15	trans-1,2-Dichloroethene	***** NOT FOUND *****		
16	MTBE	***** NOT FOUND *****		
17	Allyl chloride	***** NOT FOUND *****		
18	1,1-Dichloroethane	***** NOT FOUND *****		
19	Carbon disulfide	***** NOT FOUND *****		
20	Chloroprene	***** NOT FOUND *****		
21	Methacrylonitrile	***** NOT FOUND *****		
22	cis-1,2-Dichloroethene	***** NOT FOUND *****		
23	Methyl acrylate	***** NOT FOUND *****		
24	Chloroform	***** NOT FOUND *****		
25	2,2-Dichloropropane	***** NOT FOUND *****		
26	Dibromofluoromethane-surrogate	2.89	92.7 ppb	
27	Isobutyl alcohol	***** NOT FOUND *****		
28	Tetrahydrofuran	***** NOT FOUND *****		
29	1,1,1-Trichloroethane	***** NOT FOUND *****		
30	1,2-Dichloroethane	***** NOT FOUND *****		
31	1,1-Dichloropropene	***** NOT FOUND *****		
32	Benzene	***** NOT FOUND *****		

93 %R

33	Carbon Tetrachloride	***** NOT FOUND *****		
34	2-Chloroethanol	***** NOT FOUND *****		
35	1,4-Difluorobenzene	3.40	100.0 ppb	ISTD2
36	1,2-Dichloropropane	***** NOT FOUND *****		
37	Trichloroethene	***** NOT FOUND *****		
38	Dibromomethane	***** NOT FOUND *****		
39	Bromodichloromethane	***** NOT FOUND *****		
40	Methyl methacrylate	***** NOT FOUND *****		
41	1,4-Dioxane	***** NOT FOUND *****		
42	cis-1,3-Dichloropropene	***** NOT FOUND *****		
43	2-Nitropropane	***** NOT FOUND *****		
44	trans-1,3-Dichloropropene	***** NOT FOUND *****		
45	Toluene	***** NOT FOUND *****		
46	Toluene-d8-surrogate	4.89	103.0 ppb	103%
47	1,1,2-Trichloroethane	***** NOT FOUND *****		
48	1,3-Dichloropropane	***** NOT FOUND *****		
49	Ethylmethacrylate	***** NOT FOUND *****		
50	Dibromochloromethane	***** NOT FOUND *****		
51	1,2-Dibromomethane	***** NOT FOUND *****		
52	Tetrachloroethene	***** NOT FOUND *****		
53	Chlorobenzene	***** NOT FOUND *****		
54	Chlorobenzene-d5	7.54	100.0 ppb	ISTD3
55	1,1,1,2-Tetrachloroethane	***** NOT FOUND *****		
56	Ethylbenzene	***** NOT FOUND *****		
57	m/p-xylene	***** NOT FOUND *****		
58	Bromoform	***** NOT FOUND *****		
59	Styrene	***** NOT FOUND *****		
60	o-Xylene	***** NOT FOUND *****		
61	cis-1,4-Dichloro-2-butene	***** NOT FOUND *****		
62	trans-1,4-Dichloro-2-butene	***** NOT FOUND *****		
63	Bromofluorobenzene-surrogate	10.45	96.8 ppb	97%
64	1,1,2,2-Tetrachloroethane	***** NOT FOUND *****		
65	1,2,3-Trichloropropane	***** NOT FOUND *****		
66	Isopropylbenzene	***** NOT FOUND *****		
67	Bromobenzene	***** NOT FOUND *****		
68	2-Chlorotoluene	***** NOT FOUND *****		
69	n-Propylbenzene	***** NOT FOUND *****		
70	4-Chlorotoluene	***** NOT FOUND *****		
71	1,3,5-Trimethylbenzene	***** NOT FOUND *****		
72	Pentachloroethane	***** NOT FOUND *****		
73	tert-Butylbenzene	***** NOT FOUND *****		
74	1,2,4-Trimethylbenzene	***** NOT FOUND *****		
75	1,3-Dichlorobenzene	***** NOT FOUND *****		
76	1,4-Dichlorobenzene-d4	13.35	100.0 ppb	ISTD4
77	sec-Butylbenzene	***** NOT FOUND *****		
78	1,4-Dichlorobenzene	***** NOT FOUND *****		
79	p-iso-Propyltoluene	***** NOT FOUND *****		
80	1,2-Dichlorobenzene	***** NOT FOUND *****		
81	n-Butylbenzene	***** NOT FOUND *****		
82	1,2-Bromo-3-chloropropane	***** NOT FOUND *****		
83	Nitobenzene	***** NOT FOUND *****		
84	1,2,4-Trichlorobenzene	***** NOT FOUND *****		
85	Naphthalene	18.40	30.5 ppb	#
86	1,2,3-Trichlorobenzene	***** NOT FOUND *****		
87	Hexachlorobutadiene	***** NOT FOUND *****		

END OF REPORT

Qualifiers Not Satisfied

Fri May 26 08:40:14 2006

ESN North Atlantic
Environmental Services Network
Direct Push Sampling In-Situ Remediation Mobile & Fixed Base Chemical Analysis Integrated Site Characterization

PO Box 6752
Scarborough, ME 04070
800-880-0533 Fax: 207-883-1912

STANDARD OPERATING PROCEDURE

USEPA SW-846 Method 8260B
Water and Soil Analysis

VOLATILE ORGANIC COMPOUNDS IN WATER AND SOIL BY GAS CHROMATOGRAPHY/MASS SPECTROMETRY (GC/MS)

Property of: ESN North Atlantic
SOP No.: ESN8260WATERSOILPT
Author: Daniel M. Twomey
Reviewed By: Robert M. Burger

Date: January 31, 2004
Revision Date: Feb 10, 2006
Project specific revisions:

1.0 SCOPE AND APPLICATION

1.1 Summary of Method 8260B Analysis of Water and Soil. This method can effectively be used for the analysis of volatile organic compounds and petroleum hydrocarbons including the early to mid portion of diesel/#2 fuel oil. For low concentration analysis (0.5 to 5 ppb) up to 25 mL of water or 10 grams of soil are concentrated on a Tekmar Purge and Trap concentration device prior to analysis. Analysis of the concentrated sample is similar for both soil and water. Medium concentration soils (25 to 50 ppb) can be analyzed after extraction or preservation with purge and trap grade methanol.

1.2 Limitations. The heaviest fraction of Diesel/#2 fuel oil is not well defined by this method (final third) especially at low concentrations. This method should not be used for water and soil when aged or weathered diesel/#2 is the target of concern.

1.3 Compound Selection. The entire 8260 list, approximately 80 compounds, a subset of site specific target compounds, non-regulated volatile target compounds, fuel component mixtures or a combination of any of the previous listed choices may be analyzed.

1.4 Method Detection Limits (MDL's) are compound dependent and vary with purging efficiency and concentration. The applicable detection range of this method is compound dependent but is approximately 0.5 to 50 μg/L or μg/kg (ppb). Determination of some structural isomers (i.e., xylenes) may be hampered by coelution.

1.5 The Reporting Limit (RL) for waters and soils for individual compounds is approximately 1.0 to 100 ppb and 100 to 1000 ppb for multi-component mixtures. RLs will be proportionately higher for samples that require dilution or reduced sample size to avoid saturation of the detector.

1.6 Restrictions - This method is restricted for use by, or under the supervision of, analysts experienced in the use of gas chromatography/mass spectroscopy for measurement of volatile organics at low ppb concentrations and skilled in the interpretation of gas chromatograms and mass spectra. Each analyst must demonstrate the ability to generate acceptable results with this method.

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1.7 Toxicity - The toxicity or carcinogenicity of chemicals used in this method has not been precisely defined. Each chemical should be treated as a potential health hazard, and exposure to these chemicals should be minimized. The laboratory is responsible for maintaining awareness of OSHA regulations regarding safe handling of chemicals used in this method

2.0 METHOD DESCRIPTION

2.1 Modified Method 8260B for the analysis of water and soil provides gas chromatographic conditions for the detection of volatile organic compounds. Water and soil samples are introduced into the gas chromatograph by the purge and trap method. A temperature program is used in the gas chromatograph to separate the organic compounds. Detection is achieved using a HP GCD with a mass selective detector

2.2 Compound Identification - Compound Identifications are obtained by analyzing standards under the same conditions used for samples and comparing their mass spectra to the spectra of authentic standards. Quantitation is accomplished by comparing the response of a major (quantitation) ion relative to an internal standard using a multi-point calibration curve.

3.0 INTERFERENCES

3.1 Contamination - Samples can be contaminated by diffusion of volatile organics (particularly chlorofluorocarbons, light hydrocarbons and methylene chloride). A trip blank prepared from analyte free water carried through sampling and subsequent storage and handling can serve as a check on such contamination.

4.0 APPARATUS AND MATERIALS

4.1Gas Chromatograph – A HP GCD capable of temperature programming; equipped with variable constant differential flow controllers, oven controller and data collection system.

4.2 Primary Column – 1 X 30 m x 0.25 mm ID HP-5 capillary column or equivalent.

4.3 Detector - Mass selective detector capable of scanning from 35 to 450 amu every second or less, using 70 volts (nominal) electron energy in the electron impact ionization mode.

4.4 Data System – The data system consist of a computer system that allows for the continuous acquisition and storage of machine readable media of mass spectra obtained throughout the duration of the chromatographic program that is interfaced to the mass spectrometer. The computer is equipped with software that allows searching the GC/MS data files for ions of a specified mass and plots their ion abundance versus time or scan number. The EPA/NIST Mass Spectral Library is available on the acquisition computer.

4.5 Concentration Device – A Tekmar 3000 Purge and Trap concentration device with vocal trap or equivalent is recommended for most applications. The Purge and Trap (P&T) device consist of a 10 or 25 cc sparger vessel, and a polymer trap (a small trap containing either carbon, tenex, silica gel, or a combination of sorbing materials).

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5.0 REAGENTS

5.1 Reagent Grade Chemicals. Unless otherwise indicated, it is intended that all reagents shall conform to the specifications of the Committee on Analytical Reagents of the American Chemical Society, where such specifications are available. Other grades may be used, provided it is first ascertained that the reagent is of sufficiently high purity to permit its use without lessening the accuracy of the determination.

5.2 Methanol, CH_3OH – Purge and Trap quality or equivalent, demonstrated to be free of analytes. Store away from other solvents.

6.0 STANDARDS

6.1 Stock Standards - Stock solutions may either be prepared from pure standard materials or purchased as certified solutions. Prepare stock standards in methanol using assayed liquids or gases, as appropriate. Because of the toxicity of some of the organohalides, primary dilutions of these materials of the toxicity should be prepared in a hood.

When compound purity is assayed to be 96% or greater, the weight may be used without correction to calculate the concentration of the stock standard. Commercially prepared stock standards may be used at any concentration if they are certified by the manufacturer or by an independent source.

6.2 Stock Chemical Standards -. This information will include at a minimum:

- Vendor name supplying standards
- Name and concentration of the standard
- Lot number of the standard
- Expiration date of the standard
- any special hazards associated with the standard

6.3 Frequency of Standard Preparation - Standards should be monitored frequently by comparison to the initial calibration. Fresh standards should be prepared if this check exceeds a 30% drift. Standards for non-gases usually need to be replaced after six months or as recommended by the standard manufacturer, unless the acceptability of the standard can be documented.

6.4 Secondary Standards - Prepare secondary dilution standards, using stock standard solutions, in methanol, as needed, that contain the compounds of interest, either singly or mixed together. The secondary dilution standards will be prepared at concentrations such that the calibration standards will bracket the working range of the analytical system. Secondary dilution standards will be stored with minimal headspace for volatiles and should be checked frequently for signs of degradation or evaporation, especially just prior to preparing calibration standards from them.

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6.5 Internal Calibration Standards – The recommended internal standards are Pentafluorobenzene, 1,4-Difluorobenzene, Chlorobenzene-d5, and 1,4-Dichlorobenzene-d4. Internal standards are normally added at the mid range of the calibration curve, 100 ppb,

6.6 Surrogate Standards -The recommended surrogates are Dibromofluoromethane, Toluene-d8, and Bromofluorobenzene. Surrogate standards are normally added throughout the range of the calibration curve, 10 to 200 ppb. Advisory control limits for waters are 50 to 150 % recovery and 30 to 200 % recovery for soils.

7.0 STANDARD CODING SYSTEM

7.1 Coding System – The standard coding system is used to trace all standards back to the vendor. All standard are required to have the appropriate code assigned when they are prepared and must be labeled accordingly. The GC code will follow the format:

FGCXXXXXXWWYZ Where:

XXXXXX is the month, day, and year that the standard was made, i.e., 071404.

WW is the page in the logbook where the standard can be found, i.e., 01 to 99.

Y is where the standard fell chronologically on the page, i.e., A through Z.

Z is the logbook number where the standard can be found, i.e., 1 through 9.

8.0 Tuning

8.1 PFTBA – Perfluoro-tert-butylamine is used by the HP-GCD for tuning of the mass selective detector. The auto tune function checks the following criteria every 24 hours:

Position of mass 69
Position of mass 219
Position of mass 414
Position of isotope mass 70
Position of isotope mass 220
Position of isotope mass 415
Ratio of mass 70 to mass 69 (0.5 - 1.6%)
Ratio of mass 220 to 219 (3.2 - 5.4%)
Ratio of mass 415 to 414 (6.8 – 11.2%)
Ratio of 219 to 69, > 15%
Ratio of 414 to 69, > 0.2%

Mass 69 precursor (<= 3%)
Mass 219 precursor (<= 6%)
Mass 414 precursor (<= 12%)

Ratio of 18 to 69 (< 20%)
Ratio of 28 to 69 (< 10%)

PO Box 6752
Scarborough, ME 04070
800-880-0533 Fax: 207-883-1912

9.0 Calibration

9.1 Multilevel Calibration - In the multilevel calibration, a minimum of three levels of standards are analyzed. The standard concentrations are selected to cover the concentration range of contaminants expected. This includes a low concentration standard at or near the reporting limit (RL). In addition, the high concentration standard needs to be within the working range of both the detector and the column. Linear regression will be used with a correlation coefficient (r) of 0.95 or greater.

9.2 Continuing Calibration – At the beginning of each day a continuing calibration check standard will be analyzed at or near the mid-level of the current calibration. The system will be considered in control if the target compounds are within 30% Difference (D) when compared to the initial calibration. Compounds that exceed these criteria will be recalibrated if the standard fails after two consecutive attempts. **Note:** See section 10.10 for monitoring of system operation.

10.0 PROCEDURE AND QUALITY CONTROL

10.1 Water Method Blank – The water method blank result must be less than the reporting limit for the analysis. The water method blank is analyzed at the beginning of each day prior to the analysis of samples. If the water method blank result is above the reporting limit for the initial reading all glassware and reagents should be checked and the associated samples reanalyzed. A water method blank consist of either 10 mL or 25 mL of analyte free water, internal standards and surrogates.

10.2 Soil Method Blank – The soil method blank result must be less than the reporting limit for the analysis. The soil method blank is analyzed at the beginning of each day prior to the analysis of samples. If the soil method blank result is above the reporting limit for the initial reading all glassware and reagents should be checked and the associated samples reanalyzed. A soil method blank consists of 10 g of analyte free sand, 10 mL of analyte free water, internal standards and surrogates.

10.3 Medium Level Soil Blank – The medium soil method blank result must be less than the reporting limit for the analysis. The medium level soil method blank is analyzed prior to the analysis of medium level extracted soil samples. If the medium level soil method blank result is above the reporting limit for the initial reading all glassware and reagents should be checked and the associated samples reanalyzed. A medium level soil method blank consists of 200 μL of MeOH used for the soil extractions, 10 mL of analyte free water, internal standards and surrogates.

10.4 Water Sample Analysis - Water samples are introduced into the gas chromatograph by purge-and-trap. 10 mL or 25 mL of water sample with both the internal standard and surrogate solutions (100 ppb) are added to the sparger and purged for 11 minutes with a stream of ultra-pure helium . The helium sweeps the volatile compounds from the sparger and carries them to the polymer trap where they are sorbed. The purging and trapping process concentrates the sample on the polymer trap.

After the sample has been purged the helium flow is redirected (the valve is switched to desorb) through the polymer trap to the GC column. The polymer trap is simultaneously flash heated to desorb all of the volatile compounds. The volatile compounds are swept through a heated transfer line with helium onto the head of the column. The initial GC temperature is set at a lower temperature causing

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the volatile compounds to condense at the head of the column where separation and elution then takes place.

10.5 Soil Sample Analysis - Soil samples are introduced into the gas chromatograph by purge-and-trap. 10 g of soil sample and 10 mL of analyte free water with both the internal standard and surrogate solutions (100 ppb) are added to the sparger and purged for 11 minutes with a stream of ultra-pure helium. The helium sweeps the volatile compounds from the sparger and carries them to the polymer trap where they are sorbed. The purging and trapping process concentrates the sample on the polymer trap.

After the sample has been purged the helium flow is redirected (the valve is switched to desorb) through the polymer trap to the GC column. The polymer trap is simultaneously flash heated to desorb all of the volatile compounds. The volatile compounds are swept through a heated transfer line with helium onto the head of the column. The initial GC temperature is set at a lower temperature causing the volatile compounds to condense at the head of the column where separation and elution then takes place.

10.6 Medium Level Soil Sample Analysis – Medium level soil samples are extracted with purge and trap grade MeOH. 10 g of soil are extracted with 10 mL of purge and trap grade MeOH (1:1 extraction). A maximum of 200 μL of MeOH extract is introduced into the gas chromatograph by purge-and-trap. 200 μL (or smaller volume if a dilution is required) of soil MeOH extract is added to 10 mL of analyte free water with both the internal standard and surrogate solutions (100 ppb) are added to the sparger and purged for 11 minutes with a stream of ultra-pure helium. The helium sweeps the volatile compounds from the sparger and carries them to the polymer trap where they are sorbed. The purging and trapping process concentrates the sample on the polymer trap.

After the sample has been purged the helium flow is redirected (the valve is switched to desorb) through the polymer trap to the GC column. The polymer trap is simultaneously flash heated to desorb all of the volatile compounds. The volatile compounds are swept through a heated transfer line with helium onto the head of the column. The initial GC temperature is set at a lower temperature causing the volatile compounds to condense at the head of the column where separation and elution then takes place.

10.7 Dilutions - If the response for a peak is off-scale, i.e., beyond the calibration range of the standards, prepare a dilution of the sample with reduced sample volume. If sample dilutions are not required (result is above an action limit undiluted) then the sample result may be flagged with an " E " to indicate that the result is estimated.

10.8 Duplicates – Duplicate analysis may be required based on project specific work plans. Duplicates should be analyzed every ten samples when required. Samples should be analyzed with similar sample volumes and the results reported.

10.9 Matrix Spikes – A matrix spike (MS) and matrix spike duplicate (MSD) are not normally analyzed but may be required based on project specific work plans. MS and MSD's should be analyzed every twenty samples when required. Matrix spike standards should be prepared from volatile organic compounds, which are representative of the compounds being investigated. At a minimum the spike should include 1,1-dichloroethene, trichloroethene, chlorobenzene, toluene, and benzene. Selected water or soil samples are spiked with the appropriate compounds either at similar concentrations expected in the samples or near the mid level of the calibration curve. MS and MSD results are reported along with the associated sample results.

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10.10 Surrogate Standards –The surrogates are Dibromofluoromethane, Toluene-d8, and Bromofluorobenzene. Surrogate standards are normally added throughout the range of the calibration curve, 10 to 200 ppb. Advisory control limits for waters are 50 to 150 % recovery and 30 to 200% recovery for soils. If two or more of the surrogates fail the sample will be reanalyzed. If the surrogates fail for the reanalysis the sample will be flagged to indicate matrix interferences. If two samples in a row fail for surrogate recoveries then a calibration check standard will be analyzed to confirm that the system is in control. If the calibration check standard fails the system will be recalibrated and the effected samples reanalyzed.

10.11 Laboratory Control Sample – A laboratory control sample (LCS) is not normally analyzed but may be required based on project specific work plans. The LCS should be analyzed every twenty samples when required. The LCS should be prepared from volatile organic compounds, which are representative of the compounds being investigated. At a minimum the LCS should include 1,1-dichloroethene, trichloroethene, chlorobenzene, toluene, and benzene. Either analyte free water of clean sand (depending on the matrix being analyzed) are spiked with the appropriate compounds either at similar concentrations expected in the samples or near the mid level of the calibration curve. The LCS advisory limits are 50 to 150 % recovery for water and 30 to 200 % recovery for soils.

11.0 REFERENCES

1. "Volatile Organic Compounds in Water by Purge-and-Trap Capillary Column Gas Chromatography with Photoionization and Electrolytic Conductivity Detectors in Series", Method 502.2, Rev. 2.0 (1989); Methods for the Determination of Organic Compounds in Drinking Water", U.S. Environmental Protection Agency, Environmental Monitoring Systems Laboratory, Cincinnati, OH, EPA/600/4-88/039, December, 1988.

2. "The Determination of Halogenated Chemicals in Water by the Purge and Trap Method", Method 502.1; U.S. Environmental Protection Agency, Environmental Monitoring and Support Laboratory: Cincinnati, OH 45268, September 1986.

3. "Volatile Aromatic and Unsaturated Organic Compounds in Water by Purge and Trap Gas Chromatography", Method 503.1; U.S. Environmental Protection Agency, Environmental Monitoring and Support Laboratory: Cincinnati, OH, September 1986.

4. Glaser, J.A., Forest, D.L., McKee, G.D., Quave, S.A., Budde, W.L. "Trace Analyses for Wastewaters", Environ. Sci. Technol., 1981, 15, 1426.

5. Bellar, T.A., Lichtenberg, J.J. "The Determination of Synthetic Organic Compounds in Water by Purge and Sequential Trapping Capillary Column Gas Chromatography", U.S. Environmental Protection Agency, Environmental Monitoring and Support Laboratory: Cincinnati, OH, 45268.

6. Chang, D.Y., Lopez, I. and Yocklovich, S. G. "Determination of Kerosene and #2 Diesel in Soil by Purge and Trap vs. extraction Procedure", Journal of Soil Contamination, 1(3): 239-251 (1992).

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Standard Sample Report Format

APPENDIX D: FIXED LABORATORY REPORTS

Analytics Environmental
Katahdin Analytical Services

analytics environmental laboratory LLC

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800-929-9906
www.analyticslab.com

June 5, 2006

Mr. Karl Kasper
Woodard & Curran, Inc.
41 Hutchins drive
Portland ME 04102

RE: **Analytical Results Case Narrative**
Analytics # 56606

Dear Mr. Kasper;

Enclosed please find the analytical results for samples submitted for the above-mentioned project. The attached Cover Page lists the sample IDs, Lab tracking numbers and collection dates for the samples included in this deliverable.

Samples were analyzed Polychlorinated Biphenyls (PCBs) by EPA Method 8082.

Unless otherwise noted in the Non-conformance Summary listed below, all of the quality control (QC) criteria including initial calibration, calibration verification, surrogate recovery, holding time and method accuracy/precision for these analyses were within acceptable limits.

This Level III data package has been assembled in the following order:

Case Narrative/Non-Conformance Summary
Sample Log Sheet - Cover Page
PCB Form 1 Data Sheet for Samples and Blanks
PCB Form 10 Confirmation Results Form (if required)
PCB Form 2 Surrogate Recoveries
PCB Form 3 MS/MSD (LCS) Recoveries
PCB Form 4 Method Blank Summary (equiv.) GC/MS Logbook Sheets
PCB Form 6 Initial Calibration Data
PCB Form 7 Continuing Calibration Check
PCB Form 8 System Monitoring Compound RT Summary
Chain of Custody (COC) Forms

QC NON CONFORMANCE SUMMARY

Sample Receipt:
No QC deviations.

PCBs by EPA Method 8082:
Sample 56606-6 (SS14) was selected for MS/MSD analysis however the data wasn't usuable due to the high concentrations of PCB 1260 detected in the sample. The laboratory control samples for the batch (L05246PASB/LD05246PASB) were in control for all analytes and results were reported without qualification.

Sincerely,
ANALYTICS Environmental Laboratory, LLC

Stephen Knollmeyer
Laboratory Director

analytics environmental laboratory LLC

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Mr. Karl Kasper
Woodard & Curran
41 Hutchins Drive
Portland ME 04102

Report Number: 56606

Revision: Rev. 0

Re:

Enclosed are the results of the analyses on your sample(s). Samples were received on 24 May 2006 and analyzed for the tests listed below. Samples were received in acceptable condition, with the exceptions noted below or on the chain of custody. The results reported herein conform to the most current NELAC standards, where applicable, unless otherwise narrated in the body of the report. Please see individual reports for specific methodologies and references.

Lab Number	Sample Date	Station Location	Analysis	Comments
56606-1	05/23/06	SS01	EPA 8082 (PCBs only)	
56606-2	05/23/06	SS02	EPA 8082 (PCBs only)	
56606-3	05/23/06	SS03	EPA 8082 (PCBs only)	
56606-4	05/23/06	SS04	EPA 8082 (PCBs only)	
56606-5	05/23/06	SS13	EPA 8082 (PCBs only)	
56606-6	05/23/06	SS14	Electronic Data Deliverable	
	05/23/06	SS14	EPA 8082 (PCBs only)	

Sample Receipt Exceptions: None

Analytics Environmental Laboratory is certified by the states of New Hampshire, Maine, Massachusetts, Connecticut, Rhode Island, North Carolina, New York, Virginia, Pennsylvania and is validated by the U.S. Army Corps of Engineers (MRD) and U.S. Navy (NFESC). A list of actual certified parameters is available upon request.

If you have any further question on the analytical methods or these results, do not hesitate to call.

Authorized signature _____
Stephen L. Knollmeyer Lab. Director

Date 05/25/06

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analytics environmental laboratory LLC

EXTRACTION
LOGBOOK PAGES

✓ AK 5.24.06

Transfer Date: 05/24/06

Initials: SC

EXTRACT TRANSFER SHEET

Organochlorine PST

Sample	Analysis	Extraction Date	Matrix	Initial Volume(mL)	Initial Weight(g)	Final Volume(mL)	% Solids	Extraction Comments
56606-1	8082ASE	05/24/06	Solid		15.06	10.0		
56606-2	8082ASE	05/24/06	Solid		15.19	10.0		
56606-4	8082ASE	05/24/06	Solid		15.15	10.0		
56606-5	8082ASE	05/24/06	Solid		15.70	10.0		
56606-6	8082ASE	05/24/06	Solid		15.62	10.0		
56606-3	8082ASE	05/24/06	Solid		15.85	10.0		

✓RM
5/25/06

Pesticide and PCB Solid and Waste Preparation Logbook

Method References SW-846 3550B Sonication
SW-846 3545 Accelerated Solvent Extraction (ASE)
SW-846 3580 Concentrated Waste Dilution

Date: 5-24-06 Time: 3:30 QC Batch ID : Q-P-052406-2-SC
Initials: SC Prep. Batch ID P-P-052406-2-SC
Reviewed by: AK 5-24-06 ✓

Sample #	Matrix	Weight (g)	mL Surr (1)	mL Spike (2)	Final Vol Hexane	Cook Down Date	Vessel #	CleanUp
B05246PAS	Na2SO4	15.00	1.0	—	10mL	5/24/06	31203	A/C
L05246PASB	SiSand	15.93	1.0	1.0		AK	31159	
LD05246PASB	↓	15.61	1.0	1.0			32842	
56606-1	Soil	15.06	1.0	—			32767	
-2		15.19	1.0	—			39290	
-3		15.85	1.0	—			38279	
-4		15.15	1.0	—			31181	
-5		15.70	1.0	—			38368	
-6		15.62	1.0	—			30416	
-6.MS		15.59	1.0	1.0			38907	
-6.MSD		15.63	1.0	1.0			36776	

5/24/06 SC

Hexane Lot Number: 46060
Si Sand Lot Number: 074617 (1)
Na_2SO_4 Lot Number: 45251602
1:1 hexane and Acetone /
Diatomaceous earth /
Conc H2SO4 B41041
Florosil Cartridge /

	Lot #	Exp Date
PP Surrogate (1.0 ug/mL):	P4012	11-10-06
(2) PCB MS (4.0 ug/mL):	P3992	9-21-06
Pest Spike (500 ug/L):	—	—

03/18/05 Rev 1

[12]

analytics environmental laboratory LLC

PCB
DATA SUMMARIES

analytics environmental laboratory LLC

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603-436-5111 Fax 603-430-2151
800-929-9906

Mr. Karl Kasper
Woodard & Curran
41 Hutchins Drive
Portland ME 04102

June 5, 2006

SAMPLE DATA

CLIENT SAMPLE ID

Project Name:

Project Number:

Field Sample ID: LABQC

Lab Sample ID: B05246PAS
Matrix: Solid
Percent Solid: 100
Dilution Factor: 1.0
Collection Date:
Lab Receipt Date:
Extraction Date: 05/24/06
Analysis Date: 05/24/06

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit µg/kg	Results µg/kg
PCB-1016	15	U
PCB-1221	15	U
PCB-1232	15	U
PCB-1242	15	U
PCB-1248	15	U
PCB-1254	15	U
PCB-1260	15	U

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	84	%
Decachlorobiphenyl	68	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample was analyzed according to "Test Methods for Evaluating Solid Waste, SW-846 Method 8081."

COMMENTS: Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3545. Results are expressed on a dry weight basis.

PCB Report

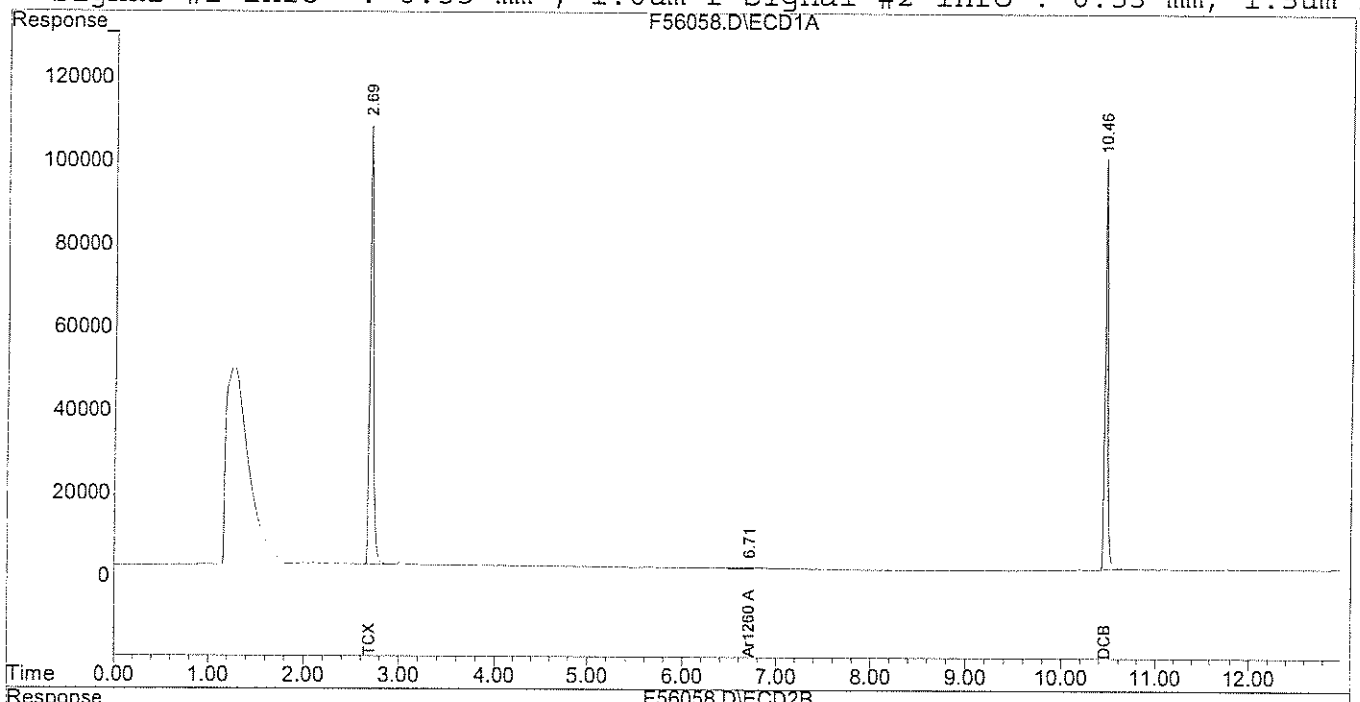
Authorized signature Melissa Gill

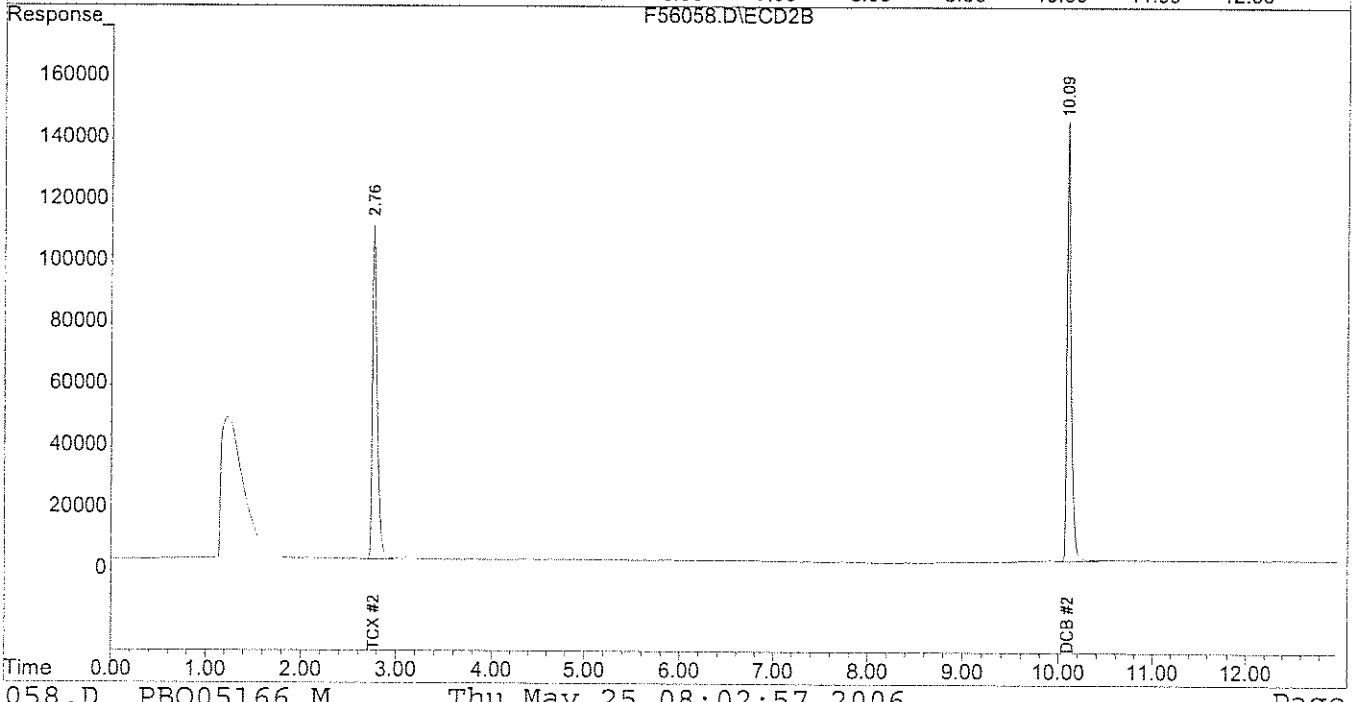
Quantitation Report (Not Reviewed)

Signal #1 : C:\HPCHEM\2\DATA\052406-F\F56058.D\ECD1A.CH Vial: 2
Signal #2 : C:\HPCHEM\2\DATA\052406-F\F56058.D\ECD2B.CH
Acq On : 24 May 2006 5:13 pm Operator:
Sample : B05246PAS, A/C Inst : Instrumen
Misc : SOIL Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 25 8:02 2006 Quant Results File: PBQ05166.RES

Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 10:06:32 2006
Response via : Multiple Level Calibration
DataAcq Meth : PSTQUICK.M

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film





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Mr. Karl Kasper
Woodard & Curran
41 Hutchins Drive
Portland ME 04102

June 5, 2006

CLIENT SAMPLE ID

Project Name:

Project Number:

Field Sample ID: LABQC

SAMPLE DATA

Lab Sample ID: B05246PAS RR
Matrix: Solid
Percent Solid: 100
Dilution Factor: 1.0
Collection Date:
Lab Receipt Date:
Extraction Date: 05/24/06
Analysis Date: 05/25/06

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit µg/kg	Results µg/kg
PCB-1016	15	U
PCB-1221	15	U
PCB-1232	15	U
PCB-1242	15	U
PCB-1248	15	U
PCB-1254	15	U
PCB-1260	15	U

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	102	%
Decachlorobiphenyl	85	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample was analyzed according to "Test Methods for Evaluating Solid Waste, SW-846 Method 8081."

COMMENTS: Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3545. Results are expressed on a dry weight basis.

PCB Report

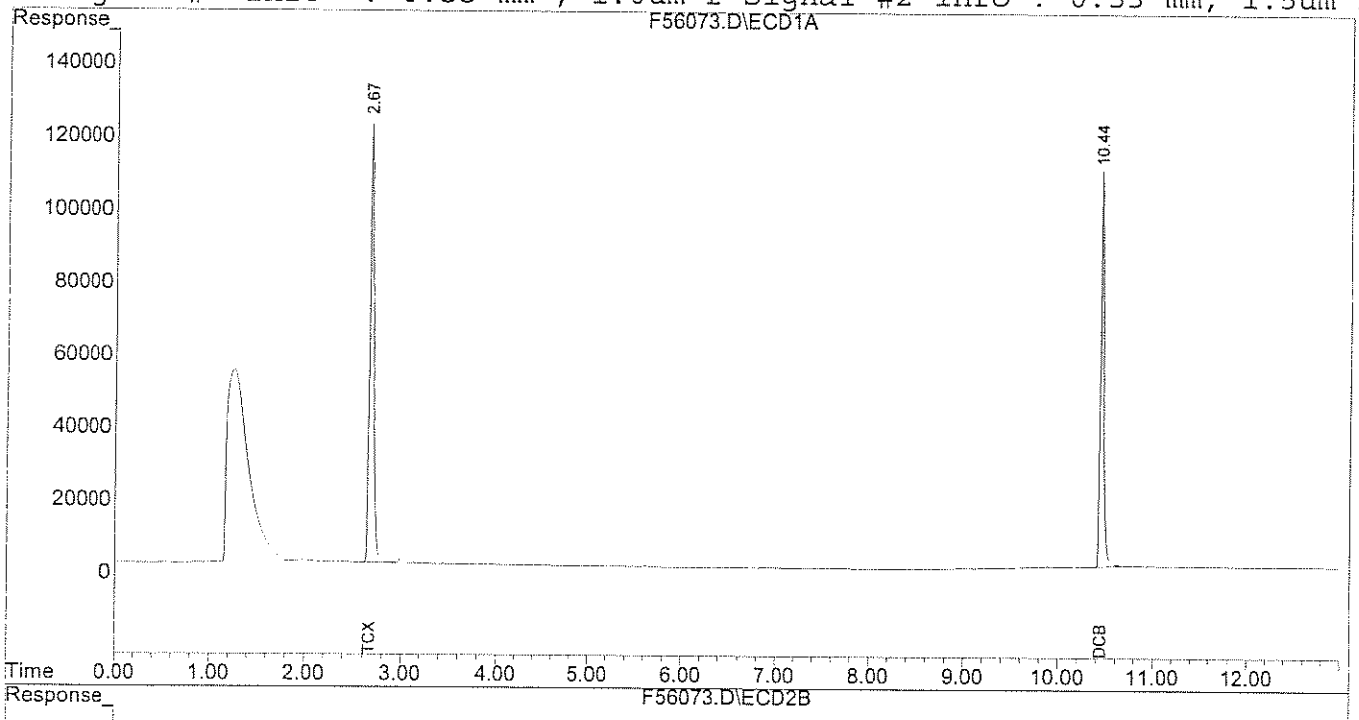
Authorized signature

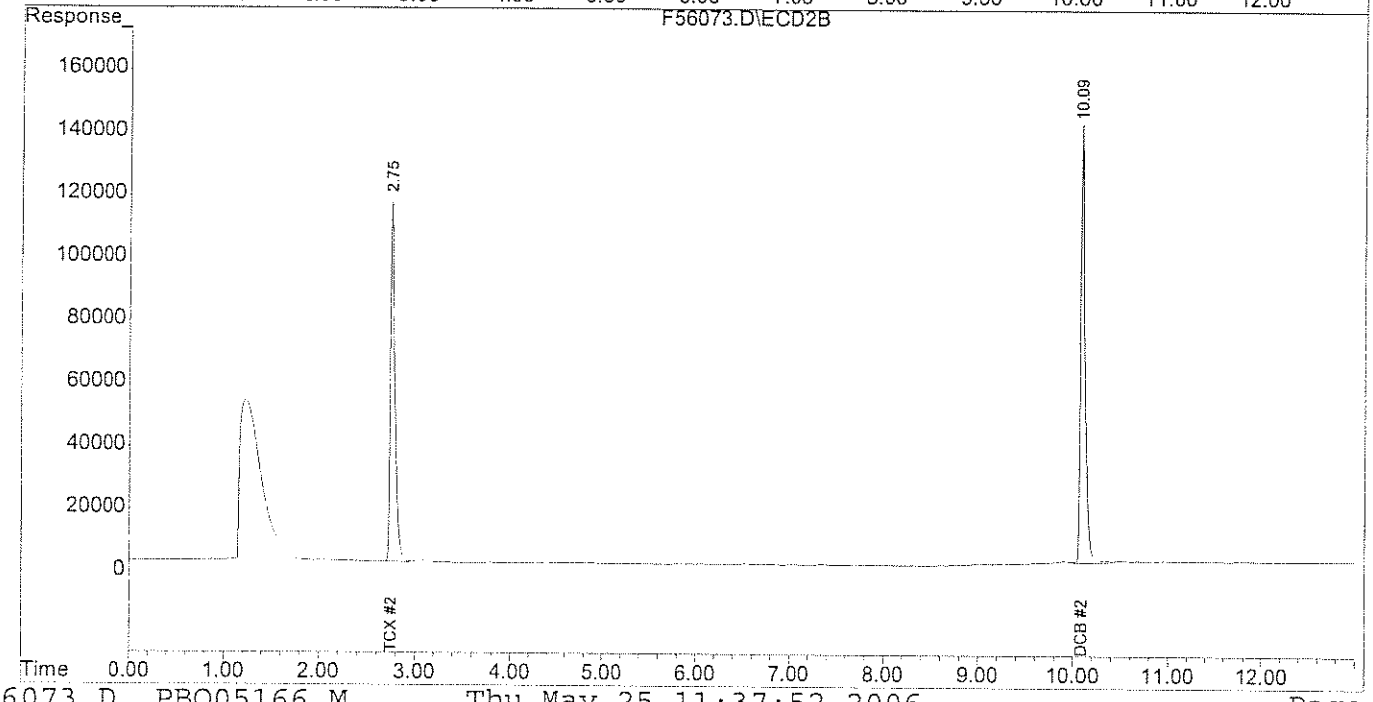
Quantitation Report (Not Reviewed)

Signal #1 : C:\HPCHEM\2\DATA\052506-F\F56073.D\ECD1A.CH Vial: 2
Signal #2 : C:\HPCHEM\2\DATA\052506-F\F56073.D\ECD2B.CH
Acq On : 25 May 2006 11:01 am Operator:
Sample : B05246PAS, A/C, RR Inst : Instrumen
Misc : SOIL Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 25 11:37 2006 Quant Results File: PBQ05166.RES

Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 10:06:32 2006
Response via : Multiple Level Calibration
DataAcq Meth : PSTQUICK.M

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film





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Mr. Karl Kasper
Woodard & Curran
41 Hutchins Drive
Portland ME 04102

May 25, 2006

SAMPLE DATA

CLIENT SAMPLE ID

Project Name:

Project Number:

Field Sample ID: SS01

Lab Sample ID:	56606-1
Matrix:	Solid
Percent Solid:	91
Dilution Factor:	1.1
Collection Date:	05/23/06
Lab Receipt Date:	05/24/06
Extraction Date:	05/24/06
Analysis Date:	05/24/06

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit µg/kg	Results µg/kg
PCB-1016	17	U
PCB-1221	17	U
PCB-1232	17	U
PCB-1242	17	U
PCB-1248	17	U
PCB-1254	17	U
PCB-1260	17	**214**

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	101	%
Decachlorobiphenyl	99	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082.

COMMENTS: Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3545. Results are expressed on a dry weight basis.

PCB Report

Authorized signature

analytics environmental laboratory LLC

PCB CONFIRMATION RESULTS

SDG: 56606

Project No: Site ID: Location: ME

Instrument ID: F GC Columns: RTX-CLPesticidesI ID: 0.32(mm)

RTX-CLPesticidesII

Sample ID: 56606-1

COMPOUND	RESPONSE #1 ug/kg	RESPONSE #2 ug/kg	DIFFERENCE %	#	% DIFFERENCE Acceptance Limit
PCB 1260	190	214	12		40

Column to be used to flag % Difference with an asterisk

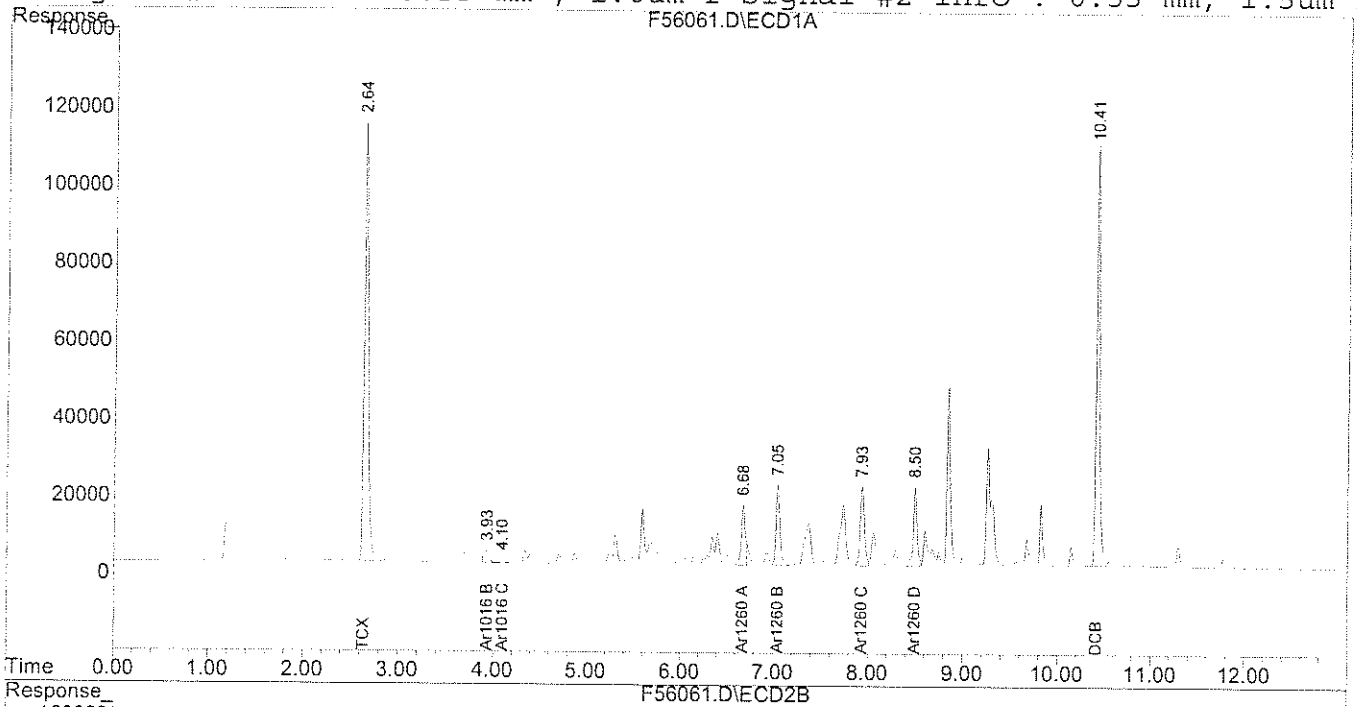
* Values outside of QC limits

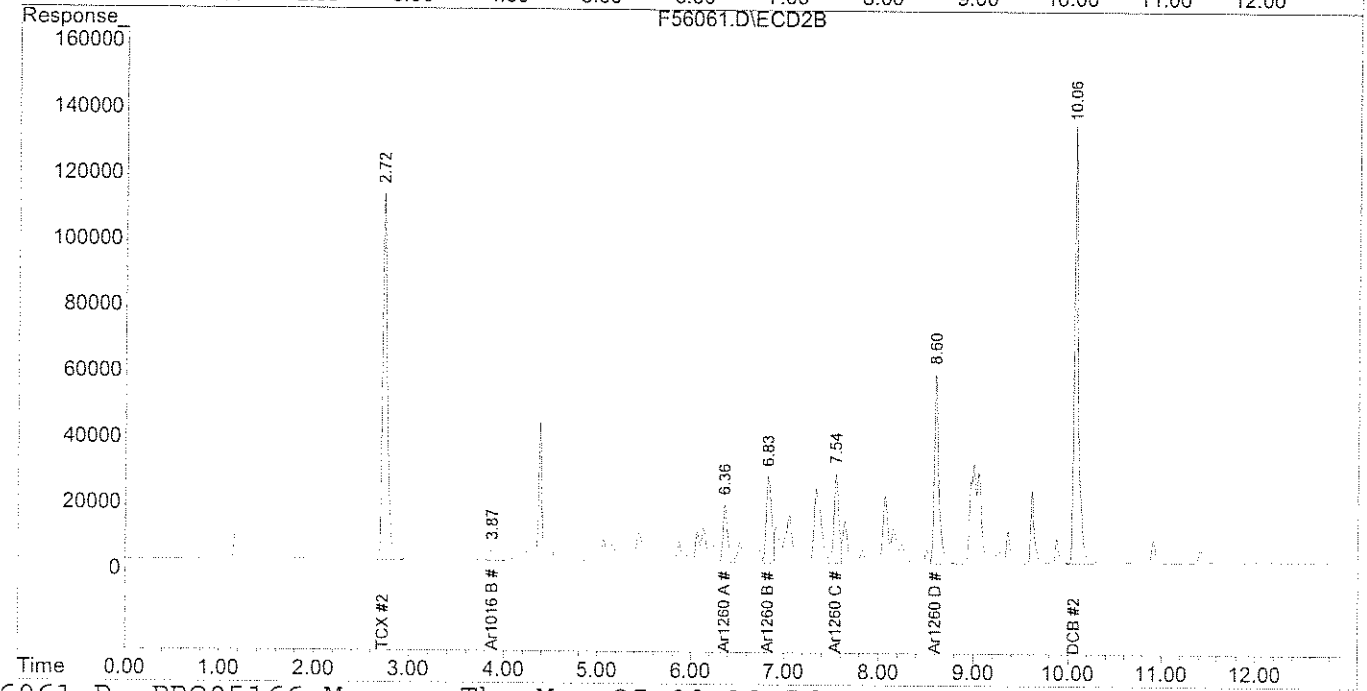
Quantitation Report (QT Reviewed)

Signal #1 : C:\HPCHEM\2\DATA\052406-F\F56061.D\ECD1A.CH Vial: 5
Signal #2 : C:\HPCHEM\2\DATA\052406-F\F56061.D\ECD2B.CH
Acq On : 24 May 2006 5:57 pm Operator:
Sample : 56606-1, A/C Inst : Instrumen
Misc : SOIL Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 25 8:21 2006 Quant Results File: PBQ05166.RES

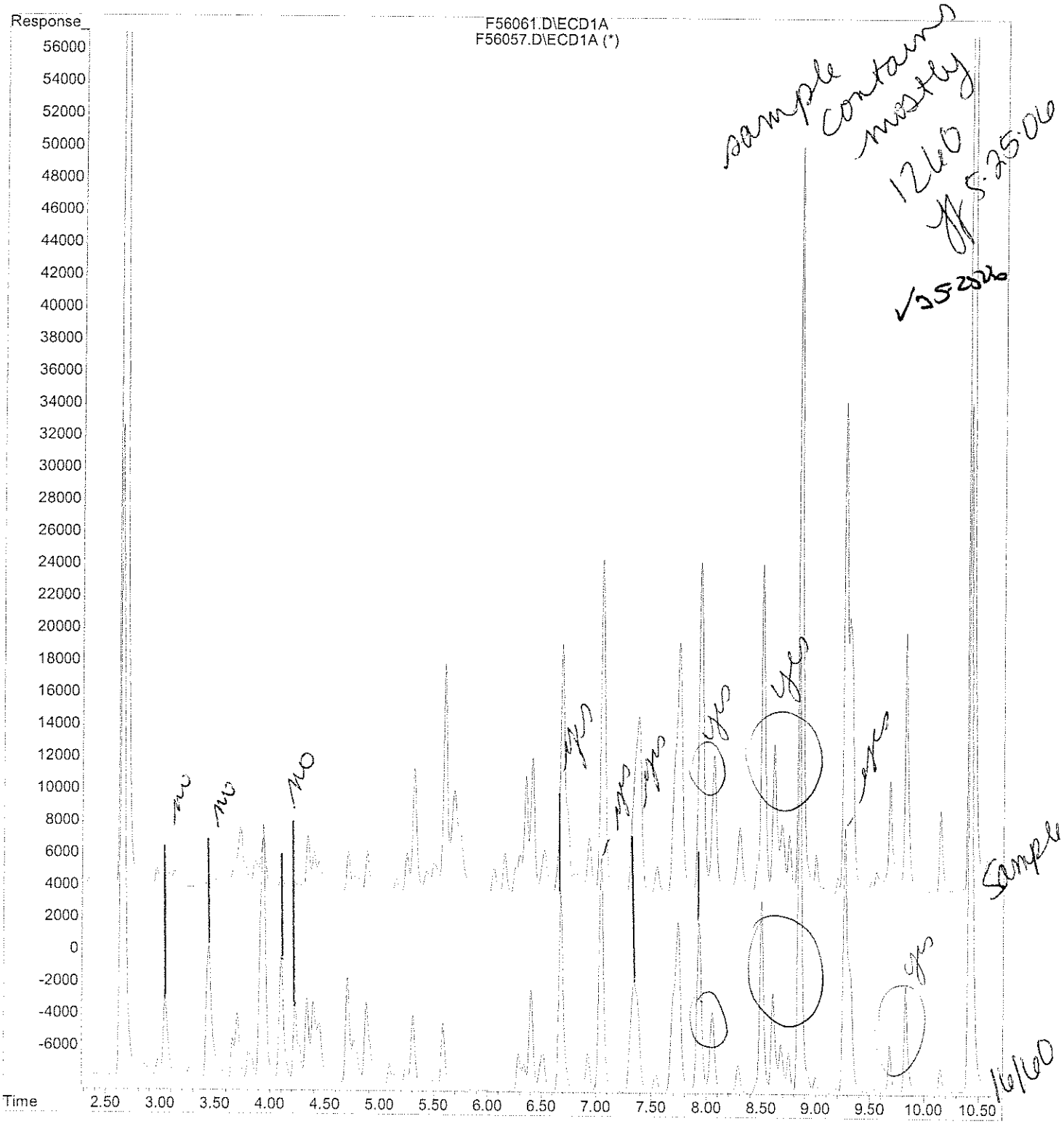
Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 10:06:32 2006
Response via : Multiple Level Calibration
DataAcq Meth : PSTQUICK.M

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film





File : C:\HPCHEM\2\DATA\052406-F\F56061.D
Operator :
Acquired : 24 May 2006 5:57 pm using AcqMethod PSTQUICK.M
Instrument : Instrumen
Sample Name: 56606-1, A/C
Misc Info : SOIL
Vial Number: 5



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Mr. Karl Kasper
Woodard & Curran
41 Hutchins Drive
Portland ME 04102

May 25, 2006

CLIENT SAMPLE ID

Project Name:

Project Number:

Field Sample ID: SS02

SAMPLE DATA

Lab Sample ID:	56606-2
Matrix:	Solid
Percent Solid:	93
Dilution Factor:	1.1
Collection Date:	05/23/06
Lab Receipt Date:	05/24/06
Extraction Date:	05/24/06
Analysis Date:	05/24/06

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit µg/kg	Results µg/kg
PCB-1016	17	U
PCB-1221	17	U
PCB-1232	17	U
PCB-1242	17	U
PCB-1248	17	U
PCB-1254	17	U
PCB-1260	17	**18**

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	91	%
Decachlorobiphenyl	78	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082.

Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3545.

COMMENTS: Results are expressed on a dry weight basis.

PCB Report

Authorized signature

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PCB CONFIRMATION RESULTS

SDG: 56606

Project No: Site ID: Location: ME

Instrument ID: F GC Columns: RTX-CLPesticidesI ID: 0.32(mm)
RTX-CLPesticidesII

Sample ID: 56606-2

COMPOUND	RESPONSE #1 ug/kg	RESPONSE #2 ug/kg	DIFFERENCE %	#	% DIFFERENCE Acceptance Limit
PCB 1260	18	13	27		40

Column to be used to flag % Difference with an asterisk

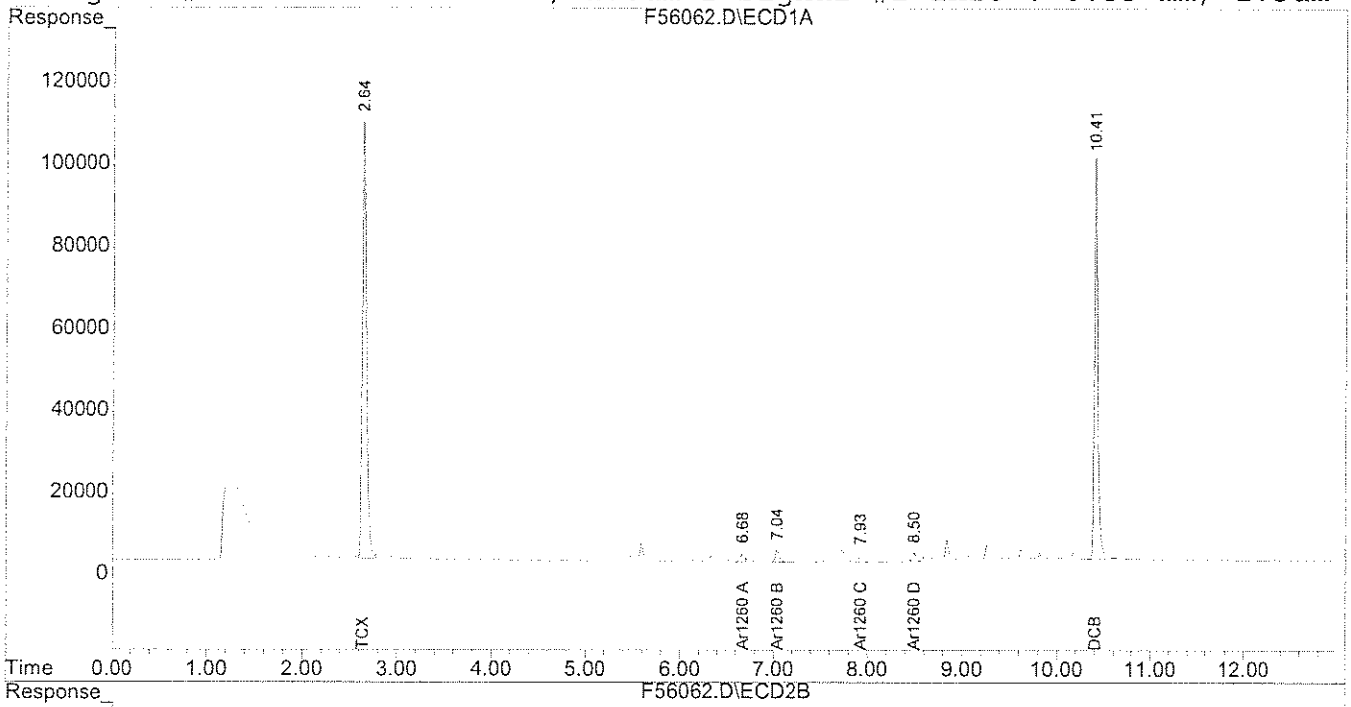
* Values outside of QC limits

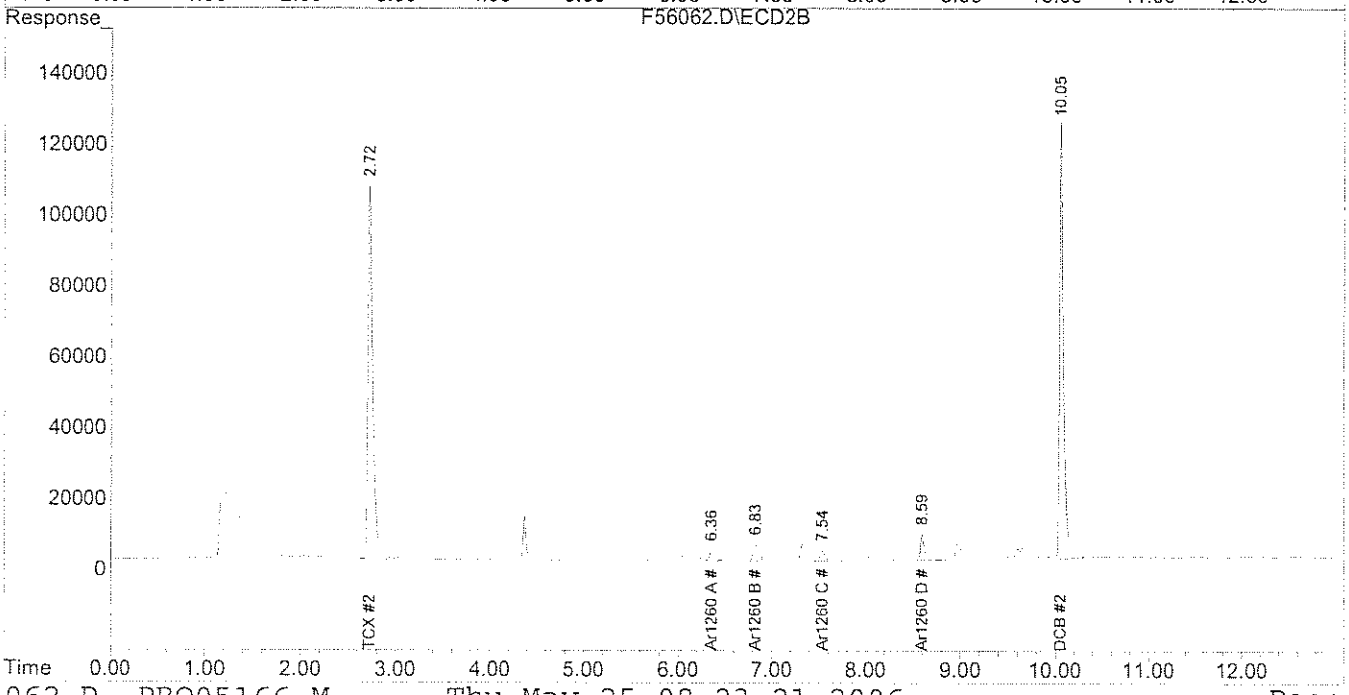
Quantitation Report (QT Reviewed)

Signal #1 : C:\HPCHEM\2\DATA\052406-F\F56062.D\ECD1A.CH Vial: 6
Signal #2 : C:\HPCHEM\2\DATA\052406-F\F56062.D\ECD2B.CH
Acq On : 24 May 2006 6:11 pm Operator:
Sample : 56606-2, A/C Inst : Instrumen
Misc : SOIL Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 25 8:23 2006 Quant Results File: PBQ05166.RES

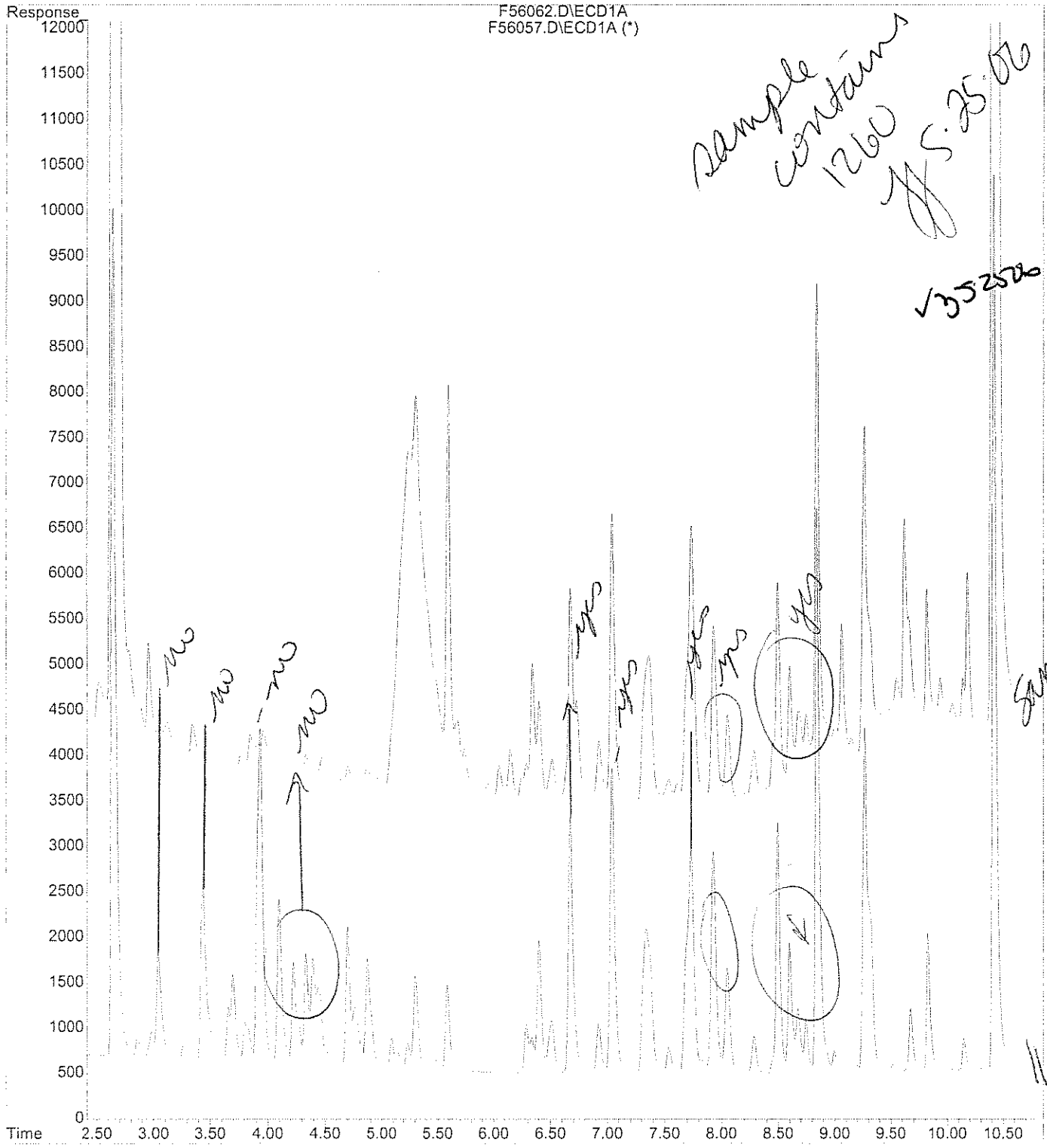
Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 10:06:32 2006
Response via : Multiple Level Calibration
DataAcq Meth : PSTQUICK.M

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film





File : C:\HPCHEM\2\DATA\052406-F\F56062.D
Operator :
Acquired : 24 May 2006 6:11 pm using AcqMethod PSTQUICK.M
Instrument : Instrumen
Sample Name: 56606-2, A/C
Misc Info : SOIL
Vial Number: 6



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Portland ME 04102

May 25, 2006

SAMPLE DATA

CLIENT SAMPLE ID

Project Name:

Project Number:

Field Sample ID: SS03

Lab Sample ID: 56606-3
Matrix: Solid
Percent Solid: 92
Dilution Factor: 1.0
Collection Date: 05/23/06
Lab Receipt Date: 05/24/06
Extraction Date: 05/24/06
Analysis Date: 05/24/06

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit µg/kg	Results µg/kg
PCB-1016	15	U
PCB-1221	15	U
PCB-1232	15	U
PCB-1242	15	U
PCB-1248	15	U
PCB-1254	15	U
PCB-1260	15	U

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	94	%
Decachlorobiphenyl	79	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082.

Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3545.

COMMENTS: Results are expressed on a dry weight basis.

PCB Report

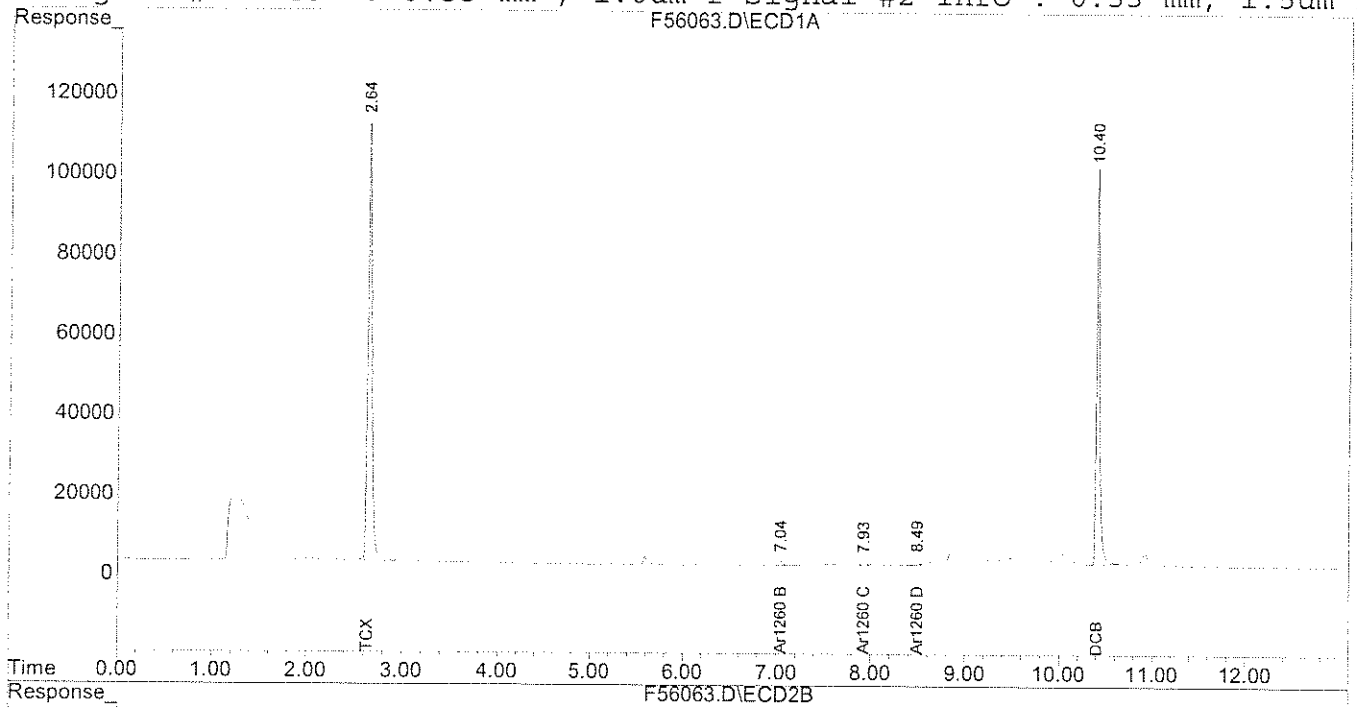
Authorized signature

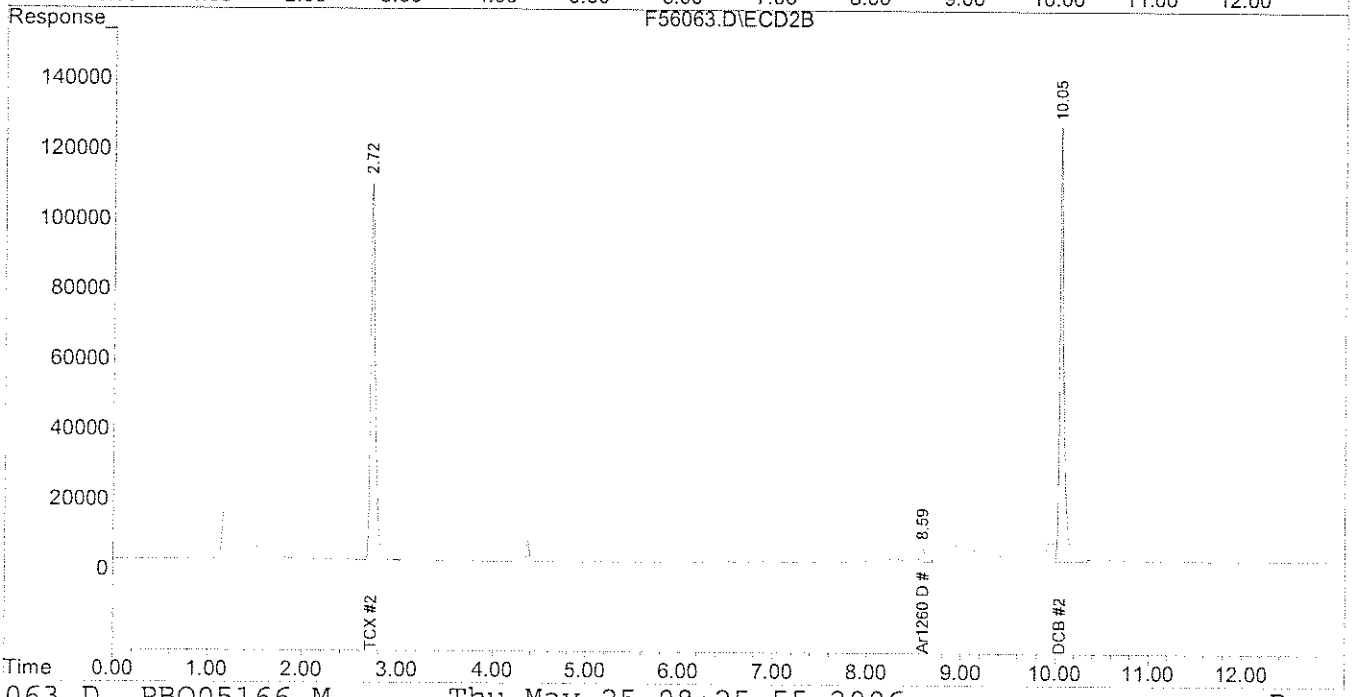
Quantitation Report (QT Reviewed)

Signal #1 : C:\HPCHEM\2\DATA\052406-F\F56063.D\ECD1A.CH Vial: 7
Signal #2 : C:\HPCHEM\2\DATA\052406-F\F56063.D\ECD2B.CH
Acq On : 24 May 2006 6:26 pm Operator:
Sample : 56606-3, A/C Inst : Instrumen
Misc : SOIL Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 25 8:25 2006 Quant Results File: PBQ05166.RES

Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 10:06:32 2006
Response via : Multiple Level Calibration
DataAcq Meth : PSTQUICK.M

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film





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May 25, 2006

SAMPLE DATA

CLIENT SAMPLE ID

Project Name:

Project Number:

Field Sample ID: SS04

Lab Sample ID:	56606-4
Matrix:	Solid
Percent Solid:	93
Dilution Factor:	1.1
Collection Date:	05/23/06
Lab Receipt Date:	05/24/06
Extraction Date:	05/24/06
Analysis Date:	05/24/06

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit µg/kg	Results µg/kg
PCB-1016	17	U
PCB-1221	17	U
PCB-1232	17	U
PCB-1242	17	U
PCB-1248	17	U
PCB-1254	17	U
PCB-1260	17	**35**

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	94	%
Decachlorobiphenyl	80	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082.

Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3545.

COMMENTS: Results are expressed on a dry weight basis.

PCB Report

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PCB CONFIRMATION RESULTS

SDG: 56606

Project No: Site ID: Location: ME

Instrument ID: F GC Columns: RTX-CLPesticidesI ID: 0.32(mm)
RTX-CLPesticidesII

Sample ID: 56606-4

COMPOUND	RESPONSE #1 ug/kg	RESPONSE #2 ug/kg	DIFFERENCE %	#	% DIFFERENCE Acceptance Limit
PCB 1260	35	32	9		40

Column to be used to flag % Difference with an asterisk

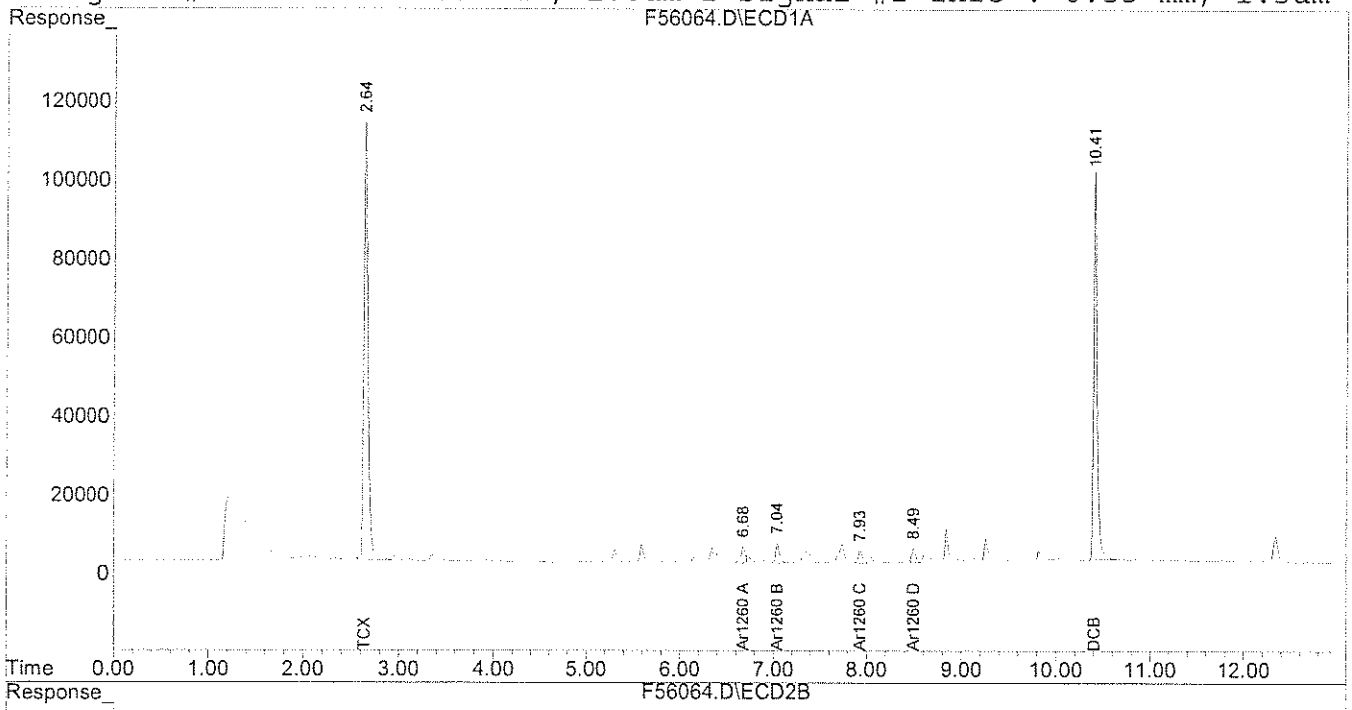
* Values outside of QC limits

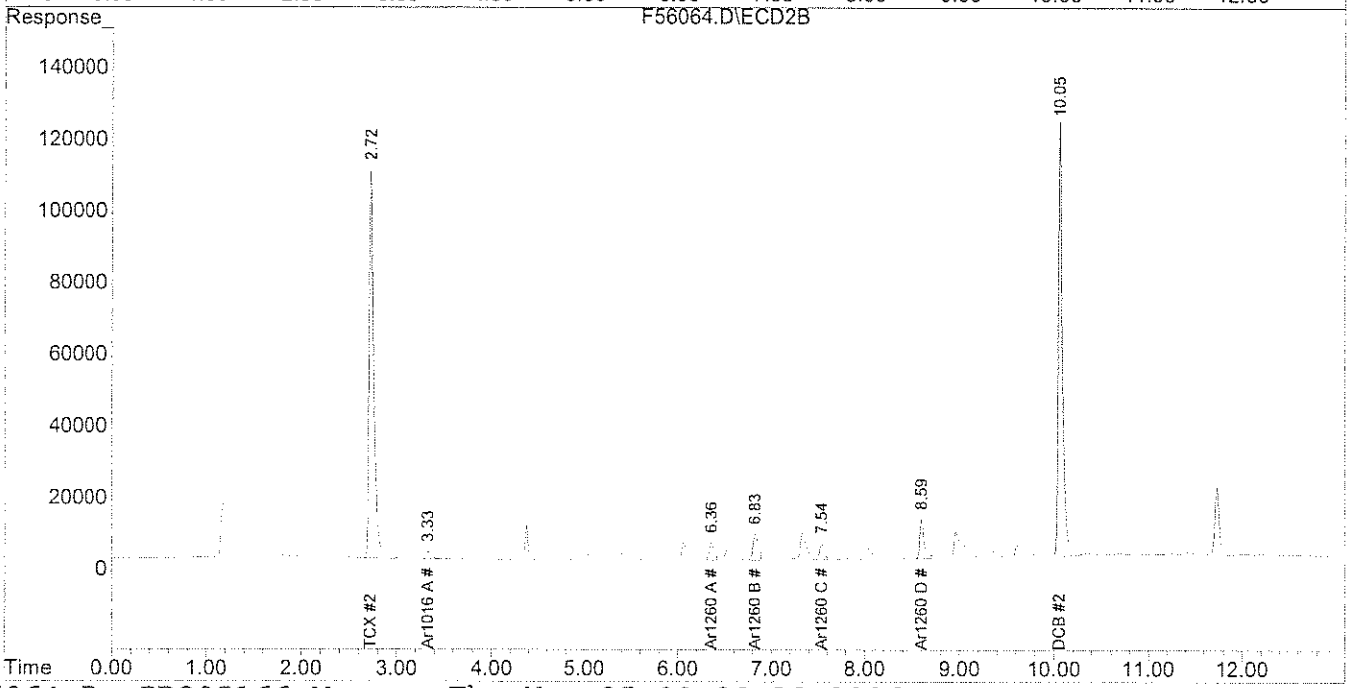
Quantitation Report (QT Reviewed)

Signal #1 : C:\HPCHEM\2\DATA\052406-F\F56064.D\ECD1A.CH Vial: 8
Signal #2 : C:\HPCHEM\2\DATA\052406-F\F56064.D\ECD2B.CH
Acq On : 24 May 2006 6:41 pm Operator:
Sample : 56606-4, A/C Inst : Instrumen
Misc : SOIL Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 25 8:28 2006 Quant Results File: PBQ05166.RES

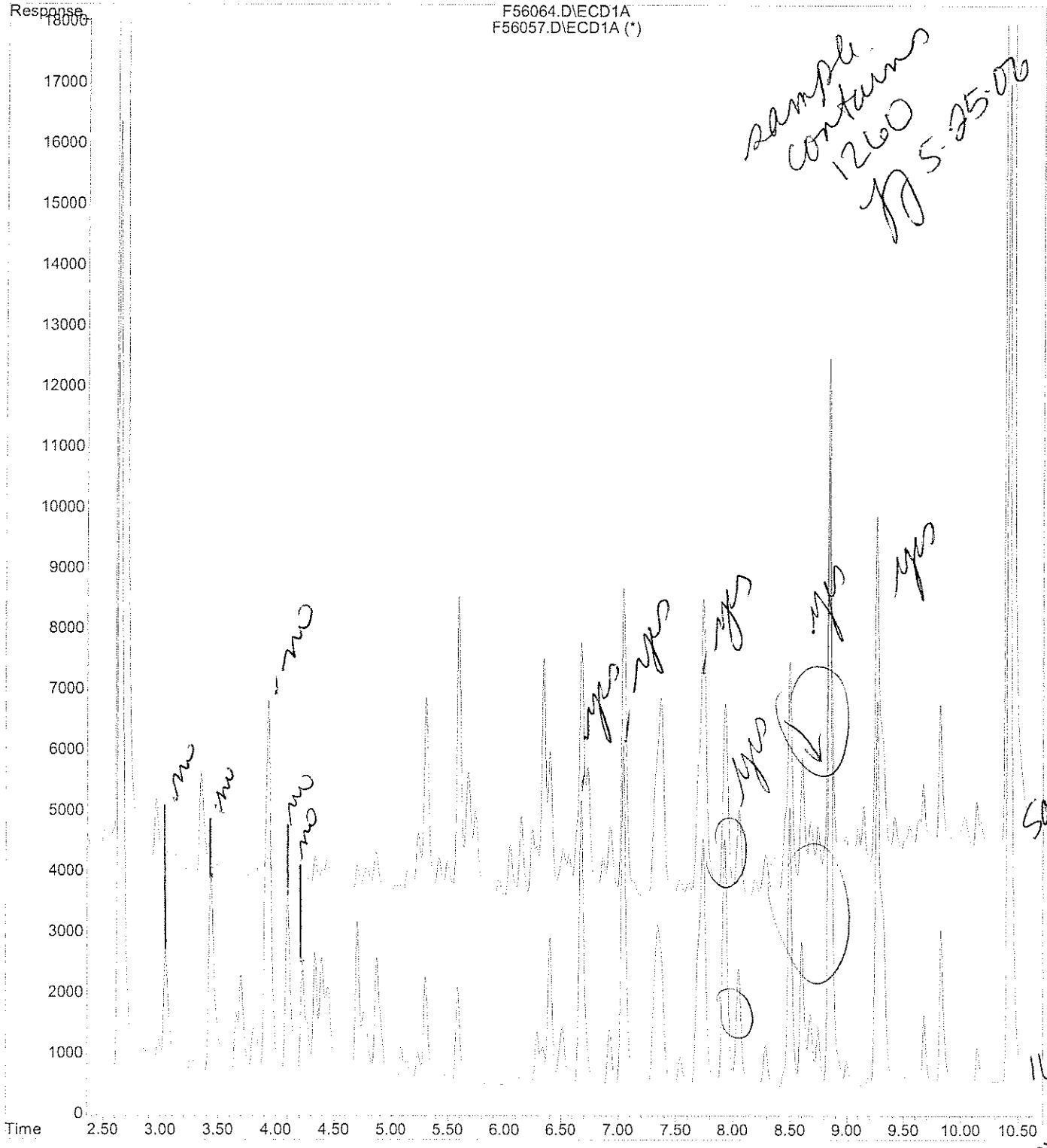
Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 10:06:32 2006
Response via : Multiple Level Calibration
DataAcq Meth : PSTQUICK.M

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film





File : C:\HPCHEM\2\DATA\052406-F\F56064.D
Operator :
Acquired : 24 May 2006 6:41 pm using AcqMethod PSTQUICK.M
Instrument : Instrumen
Sample Name: 56606-4, A/C
Misc Info : SOIL
Vial Number: 8



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Portland ME 04102

May 25, 2006

CLIENT SAMPLE ID

Project Name:

Project Number:

Field Sample ID: SS13

SAMPLE DATA

Lab Sample ID:	56606-5
Matrix:	Solid
Percent Solid:	94
Dilution Factor:	10
Collection Date:	05/23/06
Lab Receipt Date:	05/24/06
Extraction Date:	05/24/06
Analysis Date:	05/25/06

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit µg/kg	Results µg/kg
PCB-1016	150	U
PCB-1221	150	U
PCB-1232	150	U
PCB-1242	150	U
PCB-1248	150	U
PCB-1254	150	U
PCB-1260	150	**4660**

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	80	%
Decachlorobiphenyl	50	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082.

Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3545.

COMMENTS: Results are expressed on a dry weight basis.

PCB Report

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PCB CONFIRMATION RESULTS

SDG: 56606

Project No: Site ID: Location: ME

Instrument ID: F GC Columns: RTX-CLPesticidesI ID: 0.32(mm)

RTX-CLPesticidesII

Sample ID: 56606-5

COMPOUND	RESPONSE #1 ug/kg	RESPONSE #2 ug/kg	DIFFERENCE %	#	% DIFFERENCE Acceptance Limit
PCB 1260	4330	4660	7		40

Column to be used to flag % Difference with an asterisk

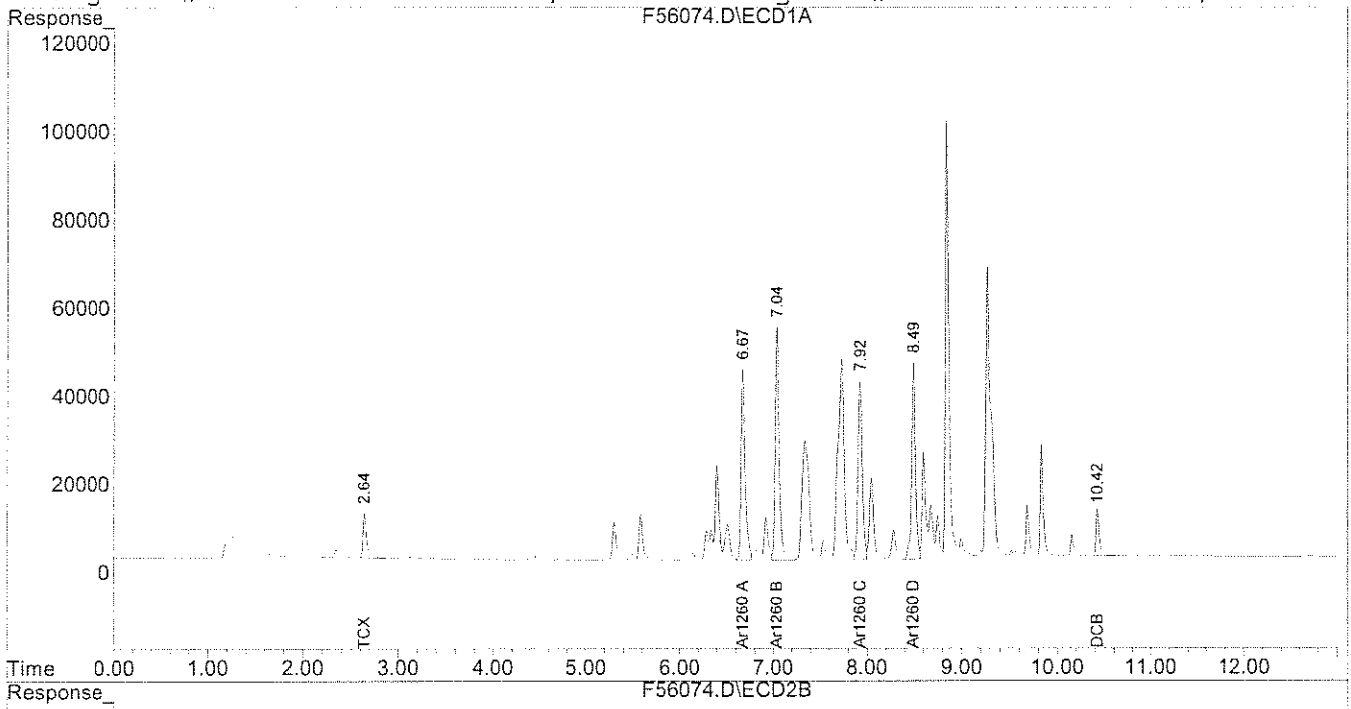
* Values outside of QC limits

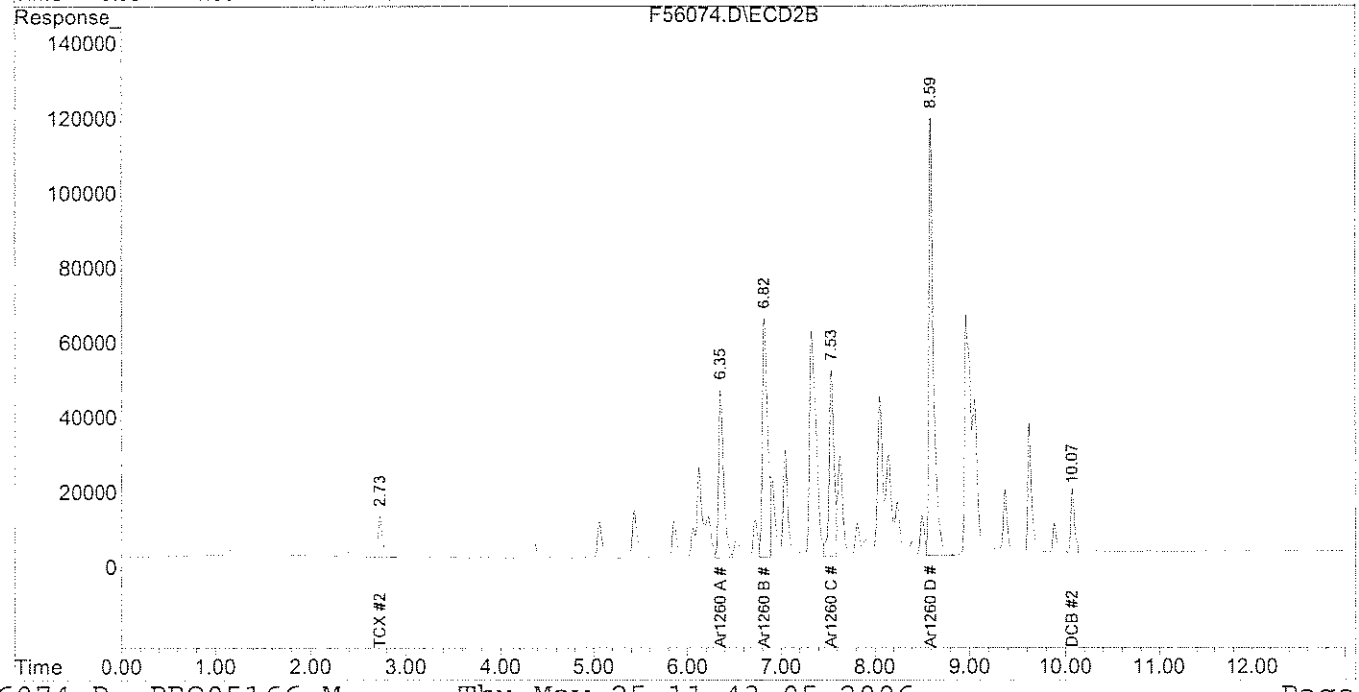
Quantitation Report (QT Reviewed)

Signal #1 : C:\HPCHEM\2\DATA\052506-F\F56074.D\ECD1A.CH Vial: 14
Signal #2 : C:\HPCHEM\2\DATA\052506-F\F56074.D\ECD2B.CH
Acq On : 25 May 2006 11:16 am Operator:
Sample : 56606-5, 1:10 Inst : Instrumen
Misc : SOIL Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 25 11:37 2006 Quant Results File: PBQ05166.RES

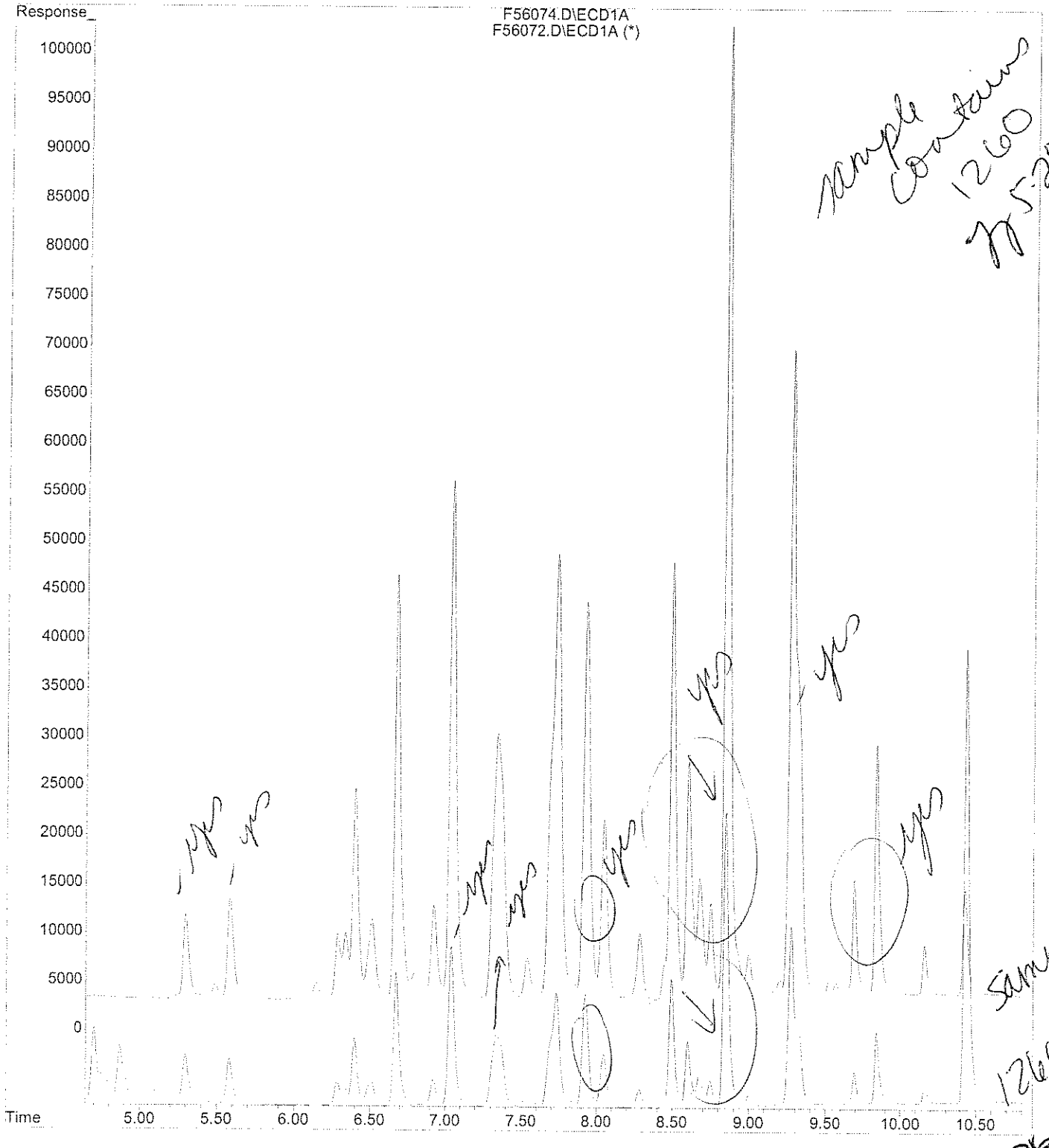
Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 10:06:32 2006
Response via : Multiple Level Calibration
DataAcq Meth : PSTQUICK.M

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film





File : C:\HPCHEM\2\DATA\052506-F\F56074.D
Operator :
Acquired : 25 May 2006 11:16 am using AcqMethod PSTQUICK.M
Instrument : Instrumen
Sample Name: 56606-5, 1:10
Misc Info : SOIL
Vial Number: 14



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Portland ME 04102

May 25, 2006

CLIENT SAMPLE ID

Project Name:

Project Number:

Field Sample ID: SS14

SAMPLE DATA

Lab Sample ID:	56606-6
Matrix:	Solid
Percent Solid:	88
Dilution Factor:	2190
Collection Date:	05/23/06
Lab Receipt Date:	05/24/06
Extraction Date:	05/24/06
Analysis Date:	05/25/06

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit µg/kg	Results µg/kg
PCB-1016	32900	U
PCB-1221	32900	U
PCB-1232	32900	U
PCB-1242	32900	U
PCB-1248	32900	U
PCB-1254	32900	U
PCB-1260	32900	**595000**

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	*	%
Decachlorobiphenyl	*	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082.

Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3545.

COMMENTS: Results are expressed on a dry weight basis. * The surrogates were diluted out.

PCB Report

Authorized signature

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PCB CONFIRMATION RESULTS

SDG: 56606

Project No: Site ID: Location: ME

Instrument ID: F GC Columns: RTX-CLPesticidesI ID: 0.32(mm)

RTX-CLPesticidesII

Sample ID: 56606-6

COMPOUND	RESPONSE #1 ug/kg	RESPONSE #2 ug/kg	DIFFERENCE %	#	% DIFFERENCE Acceptance Limit
PCB 1260	517000	595000	14		40

Column to be used to flag % Difference with an asterisk

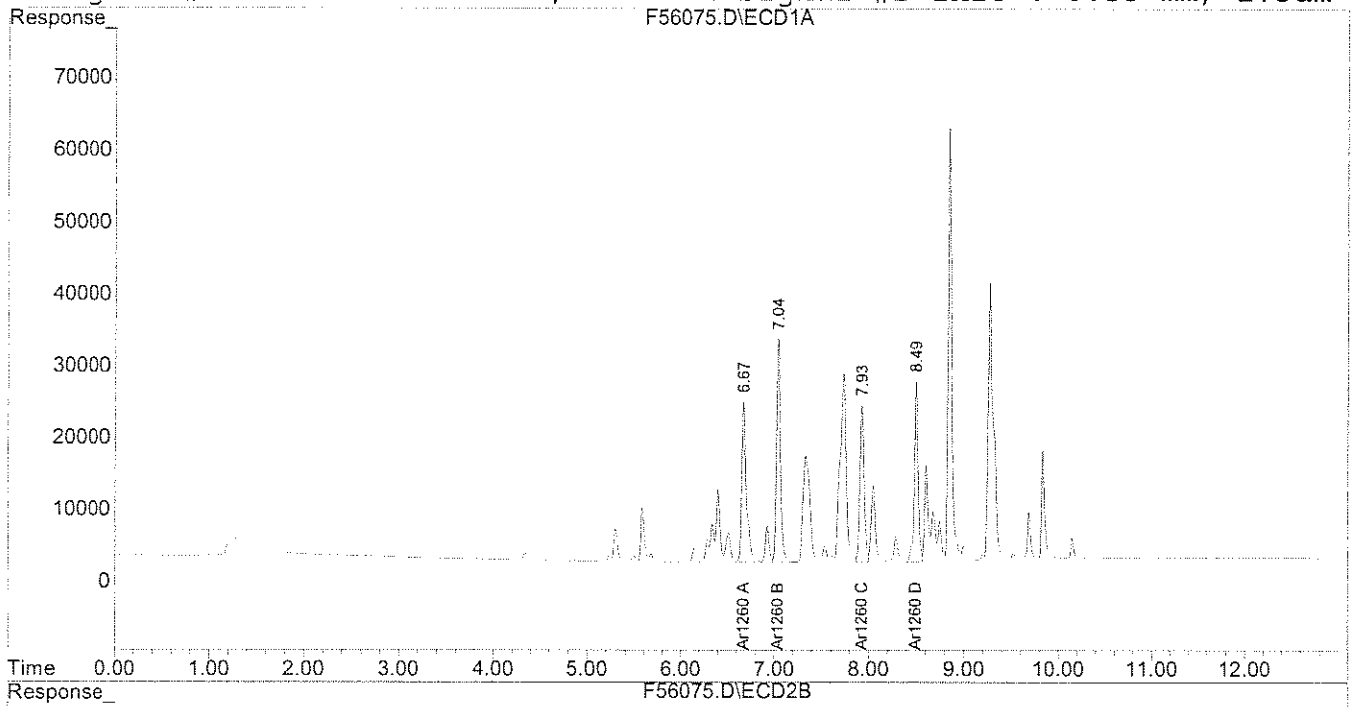
* Values outside of QC limits

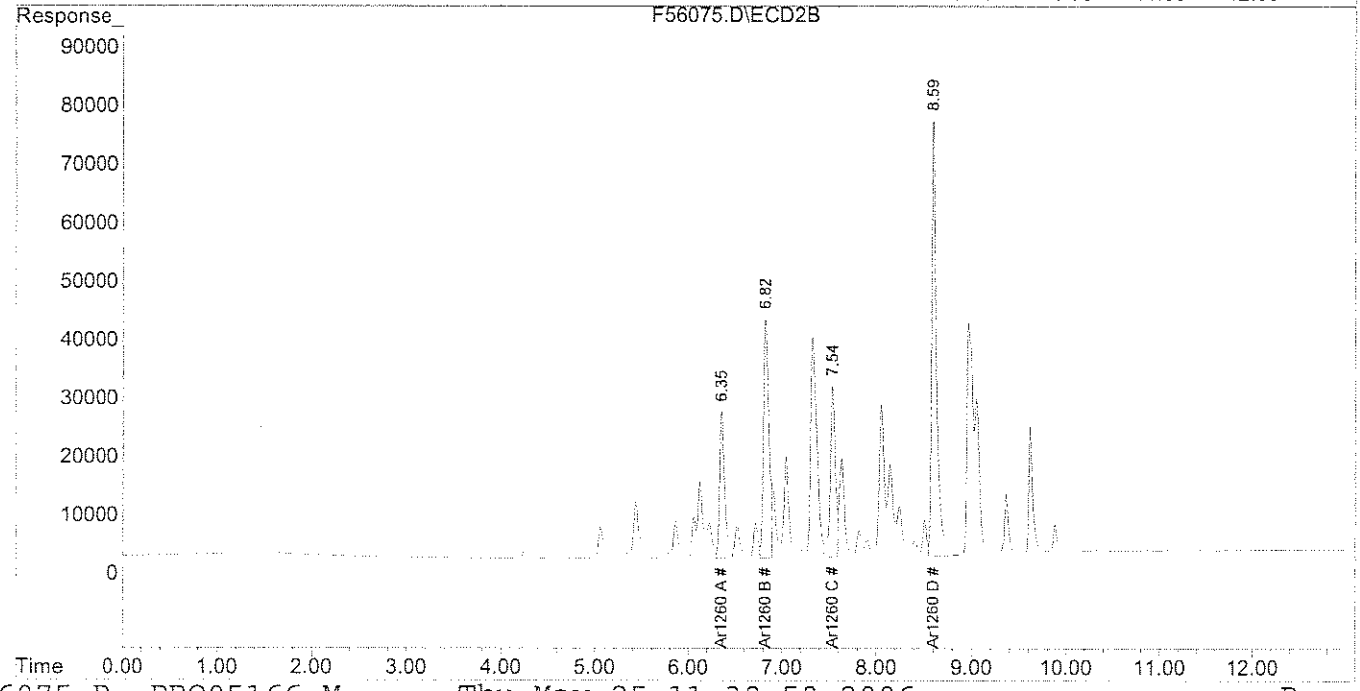
Quantitation Report (QT Reviewed)

Signal #1 : C:\HPCHEM\2\DATA\052506-F\F56075.D\ECD1A.CH Vial: 15
Signal #2 : C:\HPCHEM\2\DATA\052506-F\F56075.D\ECD2B.CH
Acq On : 25 May 2006 11:30 am Operator:
Sample : 56606-6, 1:2000 Inst : Instrumen
Misc : SOIL Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 25 11:38 2006 Quant Results File: PBQ05166.RES

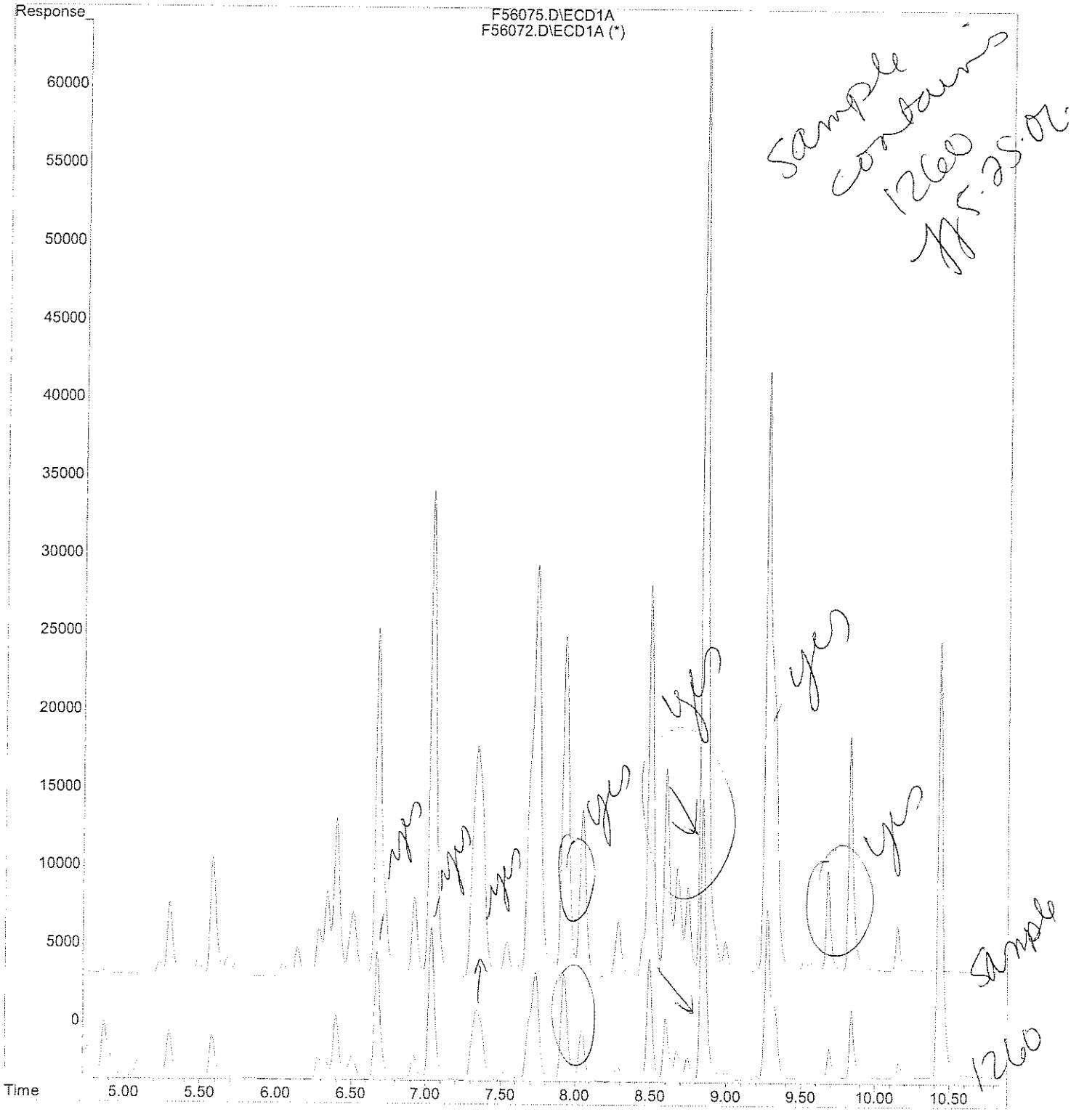
Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 10:06:32 2006
Response via : Multiple Level Calibration
DataAcq Meth : PSTQUICK.M

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film





File : C:\HPCHEM\2\DATA\052506-F\F56075.D
Operator :
Acquired : 25 May 2006 11:30 am using AcqMethod PSTQUICK.M
Instrument : Instrumen
Sample Name: 56606-6, 1:2000
Misc Info : SOIL
Vial Number: 15



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PCB
QC FORMS

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PCB SOLID
SYSTEM MONITORING COMPOUND RECOVERY

Project #: Quote #: SDG#: 56606

Site ID: Location: ME

ıstrument ID: F

GC Column: RTX-CLPesticides I ID: 0.32(mm)

RTX-CLPesticides II

		COLUMN #1			COLUMN #2			TOTAL	
	SAMPLE ID	SMC1	#	SMC2	#	SMC1	#	SMC2	#	OUT	#
01	P4016 16/60 0.01	NA		NA		NA		NA		0	
02	P4017 16/60 0.02	NA		NA		NA		NA		0	
03	P4018 16/60 0.05	NA		NA		NA		NA		0	
04	P4019 16/60 0.1	NA		NA		NA		NA		0	
05	P4020 16/60 0.25	116		109		116		121		0	
06	P4021 16/60 0.5	112		104		110		108		0	
07	P4022 16/60 1.0	NA		NA		NA		NA		0	
08	P4024 16/60 0.5 C	104		104		105		107		0	
09	P3883 AR 1221	NA		NA		NA		NA		0	
10	P3884 AR 1232	NA		NA		NA		NA		0	
11	P3885 AR 1242	NA		NA		NA		NA		0	
12	P3886 AR 1248	NA		NA		NA		NA		0	
13	P3994 AR 1254	NA		NA		NA		NA		0	
14	P3887 AR 1262	NA		NA		NA		NA		0	
15	P3888 AR 1268	NA		NA		NA		NA		0	
16	P4057 AR 1254 0.	NA		NA		NA		NA		0	
17	P4056 AR 1254 0.	NA		NA		NA		NA		0	
18	P4055 AR 1254 0.	NA		NA		NA		NA		0	
19	P4054 AR 1254 0.	NA		NA		NA		NA		0	
20	P4053 AR 1254 0.	NA		NA		NA		NA		0	
21	P4052 AR 1254 1.	NA		NA		NA		NA		0	
22	P4010 AR 1254 0.	102		103		105		108		0	
23	ZZZZ	81		17	*	87		14	*	2	
24	ZZZZ	85		21	*	95		11	*	2	
25	ZZZZ	87		15	*	95		16	*	2	
26	ZZZZ	86		14	*	94		15	*	2	
27	ZZZZ	77		81		80		81		0	
28	ZZZZ	62		33	*	87		33	*	2	
29	ZZZZ	83		10	*	96		9	*	2	
30	P4021 16/60 0.5	99		103		102		109		0	
31	P4053 AR 1254 0.	NA		NA		NA		NA		0	
32											
33											
34											
35											
36											
37											
38											
39											
40											
41											
42											

QC LIMITS

SMC1 = TCX 46-123

SMC2 = DCB 38-123

*=Values outside of protocol required QC limits

D=System Monitoring Compound diluted out

analytics environmental laboratory LLC

PCB SOLID
SYSTEM MONITORING COMPOUND RECOVERY

Project #: Quote #: SDG#: 56606

Site ID: Location: ME

ıstrument ID: F
GC Column: RTX-CLPesticides I ID: 0.32(mm)
RTX-CLPesticides II

		COLUMN #1			COLUMN #2			TOTAL	
	SAMPLE ID	SMC1	#	SMC2	#	SMC1	#	SMC2	#	OUT	#
01	P4020 16/60 0.25	98		100		104		108		0	
02	B05246PAS, A/C	84		68		99		103		0	
03	L05246PASB, A/C	98		86		107		103		0	
04	LD05246PASB, A	100		90		108		103		0	
05	56606-1, A/C	95		87		101		99		0	
06	56606-2, A/C	91		78		97		88		0	
07	56606-3, A/C	94		79		98		94		0	
08	56606-4, A/C	94		80		99		86		0	
09	ZZZZZ	81		77		85		72		0	
10	ZZZZZ	86		240	*	90		204	*	2	
11	56606-6, MS, A/C	86		126	*	91		128	*	2	
12	56606-6, MSD, A/	87		139	*	93		144	*	2	
13	P4021 16/60 0.5	98		97		104		97		0	
14											
15											
16											
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42											

QC LIMITS
SMC1 = TCX 46-123
SMC2 = DCB 38-123
*=Values outside of protocol required QC limits
D=System Monitoring Compound diluted out

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PCB SOLID
SYSTEM MONITORING COMPOUND RECOVERY

Project #: Quote #: SDG#: 56606

Site ID: Location: ME

ıstrument ID: F
GC Column: RTX-CLPesticides I ID: 0.32(mm)
RTX-CLPesticides II

		COLUMN #1		COLUMN #2		TOTAL	
	SAMPLE ID	SMC1 #	SMC2 #	SMC1 #	SMC2 #	OUT	#
01	P4020 16/60 0.25	98	88	108	114	0	
02	B05246PAS, A/C,]	102	85	107	104	0	
03	56606-5, 1:10	80	50	80	50	0	
04	56606-6, 1:2000	D	D	D	D	0	
05	P4021 16/60 0.5	101	91	104	101	0	
06							
07							
08							
09							
10							
11							
12							
13							
14							
15							
16							
17							
18							
19							
20							
21							
22							
23							
24							
25							
26							
27							
28							
29							
30							
31							
32							
33							
34							
35							
36							
37							
38							
39							
40							
41							
42							

QC LIMITS
SMC1 = TCX 46-123
SMC2 = DCB 38-123
*=Values outside of protocol required QC limits
D=System Monitoring Compound diluted out

analytics environmental laboratory LLC

PCB SOLID
LABORATORY CONTROL SPIKE/ DUPLICATE
PERCENT RECOVERY

Project #: Quote #: SDG#: 56606

Site ID: Location: ME

Instrument ID: F

GC Column: RTX-CLPesticides I ID: 0.32(mm)

RTX-CLPesticides II

Laboratory Control Spike/Duplicate Sample ID: L05246PASB A/C/ LD05246PASB A/C

COMPOUND	SPIKE ADDED (ug/kg)	SAMPLE CONCENTRATION (ug/kg)	LCS CONCENTRATION (ug/kg)	LCS RECOVERY %	#	QC LIMITS % REC
PCB 1016	267	U	244	91		67-144
PCB 1260	267	U	224	84		58-155
PCB 1016 #2	267	U	268	100		69-140
PCB 1260 #2	267	U	279	104		67-149

COMPOUND	SPIKE ADDED (ug/kg)	LCSD CONCENTRATION (ug/kg)	LCSD % REC	#	RPD	#	QC LIMITS RPD	QC LIMITS % REC.
PCB 1016	267	255	95		4		30	67-144
PCB 1260	267	238	89		6		30	58-155
PCB 1016 #2	267	273	102		2		30	69-140
PCB 1260 #2	267	283	106		2		30	67-149

Column to be used to flag recovery and RPD values with an asterisk

* Values outside of QC limits
RPD: 0 out of 4 outside limits
Spike Recovery: 0 out of 8 outside limits

Comments: _____

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PEST/PCB/HERB
Instrument F
Sequence Log Book

051606 - F

Load: gg 5.12.06
Unload: gg 5.12.06

Data File	Vial #	Sample	Method	Comments
F55857	9	P4045 Pest50		✓
58	10	P4046 Pest100		✓
59	11	P4047 Pest150		✓
60	12	P3965 Pest 0.5CC		✓
61	13	P4016 16/60 0.01	PBQ05166	✓
62	14	P4017 16/60 0.02		✓
63	15	P4018 16/60 0.05		✓
64	16	P4019 16/60 0.1		✓
65	17	P4020 16/60 0.25		✓
66	18	P4021 16/60 0.5		✓
67	19	P4022 16/60 1.0		✓
68	20	P4024 16/60 0.5CC		✓
69	21	P3883 AR1221		for overlay, single pos
70	22	P3884 AR1232		
71	23	P3885 AR 1242		
72	24	P3886 AR 1248		
73	25	P3994 AR 1254		
74	26	P3887 AR 1262		
75	27	P3888 AR 1268		
76	28	P4057 AR1254 0.02	gg 5.18.06 #S4005166	✓
77	29	P4056 AR 1254 0.05		✓
78	30	P4055 AR1254 0.1		✓
79	31	P4054 AR1254 0.25		✓
80	32	P4053 AR1254 0.5		✓
81	33	P4052 AR1254 1.0		✓
82	34	P4010 AR1254 0.5		✓
83	35	B051106PW		✓
84	36	L051106PWT		✓ methoxy ↑#1, ok#2, endrin
85	37	L051106PWB		✓ gg 5.12.06 Ketone ↑(marg.)#1 ↑#2

Standard Dilutions: 1:2- 500 uL sample +500 uL hexane
1:5- 200 uL sample +800 uL hexane
Hexane Lot : _____ 1:10- 100 uL sample + 900 uL hexane
1:20- 50 uL sample + 950 uL hexane
MTBE Lot: _____ 1:100- 10 uLsample +990 uL hexane
1:1000- 1 uLsample +1000 uL hexane

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PEST/PCB/HERB
Instrument F
Sequence Log Book

Load: YJ
Unload: JJ

052406-F.

Data File	Vial #	Sample	Method	Comments
F560506	100	Hexane		
57	1	P4D20 16160 0.25	PBQ05166	✓
58	2	B05246PAS A/C		✓
59	3	L05246PASB		✓
60	4	LD05246PASB		✓
61	5	56606-1		✓
62	6	-2		✓
63	7	-3		✓
64	8	-4		✓
65	9	-5		RR@ 1:10
66	10	-6		RR@ 1:2000
67	11	-6.MS		↑1260 hit in sample
68	12	-6.MSD		msmsD failed
69	13	P4021 16160 0.5	↓	✓

5.25.00

Standard Dilutions: 1:2- 500 uL sample +500 uL hexane
1:5- 200 uL sample +800 uL hexane
Hexane Lot : _____ 1:10- 100 uL sample + 900 uL hexane
1:20- 50 uL sample + 950 uL hexane
MTBE Lot: _____ 1:100- 10 uLsample +990 uL hexane
1:1000- 1 uLsample +1000 uL hexane

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PEST/PCB/HERB
Instrument F
Sequence Log Book

052506-F

Load: JJ
Unload: JJ

Data File	Vial #	Sample	Method	Comments
F560701	1	P4020 16|60 0.25		not eluting would peak sys. conditioning
71	100	Hexane		
72	1	P4020 16|60 0.25	PBQ051666	✓
73	2	B052466PAS A|C, RR		✓
74	14	566006-5, 1:10		✓
75	15	566006-6, 1:2000 ↓		✓
76	13	P4021 16|60 0.5		✓
77	100	Hexane		changed liner JJ 5.25.06
~~78~~	~~100~~	~~Hex~~		
78	1	P4020 16|60 0.25	PBQ051666	✓
79	2	B052566P Wipe, A|C		✓
80	3	L052566P Wipe		✓
81	4	LD052566P Wipe ↓		✓
82	5	56617-24		✓
83	6	-25		✓
84	7	-26		✓
85	8	-27		✓
86	9	-28		RR@1:10
87	10	-29		✓
88	11	-30		✓
89	12	-31		✓
90	13	-32		✓
91	14	-33		✓
92	15	-34		✓
93	16	-35		✓
94	17	-36		✓
95	18	-37 ↓		✓
96	19	P4021 16|60 0.5		✓
97	20	B052566PAS A|C		✓

Standard Dilutions: 1:2- 500 uL sample +500 uL hexane
1:5- 200 uL sample +800 uL hexane
Hexane Lot : _____ 1:10- 100 uL sample + 900 uL hexane
1:20- 50 uL sample + 950 uL hexane
MTBE Lot: _____ 1:100- 10 uLsample +990 uL hexane
1:1000- 1 uLsample +1000 uL hexane

C:\HPCHEM\2\METHODS\PBQ05166.M

Level 1 File Name	F55861.D
Level 2 File Name	F55862.D
Level 3 File Name	F55863.D
Level 4 File Name	F55864.D
Level 5 File Name	F55865.D
Level 6 File Name	F55866.D
Level 7 File Name	F55867.D

5-17-06

5-17-06

RESPONSE = X + YA + Z(A*A)

Name	R^2	R	X	Y	Z
TCX	1.000	1.000	1.04E+04	3.47E+07	0.00E-01
Ar1016 A	0.999	0.999	2.45E+04	1.65E+06	0.00E-01
Ar1016 B	0.999	1.000	3.80E+04	3.18E+06	0.00E-01
Ar1016 C	1.000	1.000	1.15E+04	1.34E+06	0.00E-01
Ar1260 A	0.999	1.000	2.75E+04	2.04E+06	0.00E-01
Ar1260 B	1.000	1.000	2.73E+04	2.45E+06	0.00E-01
Ar1260 C	1.000	1.000	8.21E+03	1.96E+06	0.00E-01
Ar1260 D	1.000	1.000	7.55E+03	1.93E+06	0.00E-01
DCB	0.999	0.999	9.77E+04	2.91E+07	0.00E-01
TCX #2	0.999	1.000	6.78E+04	3.70E+07	0.00E-01
Ar1016 A #2	0.997	0.999	4.34E+04	1.69E+06	0.00E-01
Ar1016 B #2	0.998	0.999	8.89E+04	3.83E+06	0.00E-01
Ar1016 C #2	0.997	0.999	3.37E+04	1.45E+06	0.00E-01
Ar1260 A #2	0.995	0.998	8.05E+04	2.28E+06	0.00E-01
Ar1260 B #2	0.996	0.998	1.10E+05	3.42E+06	0.00E-01
Ar1260 C #2	0.997	0.999	6.09E+04	2.54E+06	0.00E-01
Ar1260 D #2	0.999	0.999	7.76E+04	5.97E+06	0.00E-01
DCB #2	0.994	0.997	2.64E+05	3.87E+07	0.00E-01

Quantitation Report (QT Reviewed)

Signal #1 : C:\HPCHEM\2\DATA\051606-F\F55861.D\ECD1A.CH Vial: 13
Signal #2 : C:\HPCHEM\2\DATA\051606-F\F55861.D\ECD2B.CH
Acq On : 16 May 2006 7:38 pm Operator:
Sample : P4016 16/60 0.01 Inst : Instrumen
Misc : Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 17 9:56 2006 Quant Results File: PBQ05166.RES

Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 09:54:21 2006
Response via : Initial Calibration
DataAcq Meth : PSTQUICK.M

5.17.06 5.17.06

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film

Compound	RT#1	RT#2	Resp#1	Resp#2	ug/ml	ug/ml
System Monitoring Compounds						
1) S TCX	2.64	2.72	61307	77106	0.002	0.001 #
Spiked Amount 0.100	Range 46 - 122		Recovery =		2.00%#	1.00%#
9) S DCB	10.39	10.04	79033	133576	N.D. m	N.D. m
Spiked Amount 0.100	Range 26 - 86		Recovery =		0.00%#	0.00%#
Target Compounds						
2) L1 Ar1016 A	3.43	3.37	19531	27177	N.D. m	N.D.
3) L1 Ar1016 B	3.92	3.86	43320	57040	0.001	N.D. #
4) L1 Ar1016 C	4.09	4.03	15699	21986	0.004	N.D. #
Sum Ar1016 A			59019	0	0.001	N.D.
Average Ar1016 A					0.000	0.000
5) L2 Ar1260 A	6.67	6.35	30275	39717	0.002	N.D. m#
6) L2 Ar1260 B	7.03	6.81	33089	60122	0.001	N.D. m#
7) L2 Ar1260 C	7.92	7.53	22273	35850	0.006	N.D. m#
8) L2 Ar1260 D	8.49	8.59	22028	68293	0.004	N.D. #
Sum Ar1260 A			107665	0	0.013	N.D.
Average Ar1260 A					0.003	0.000

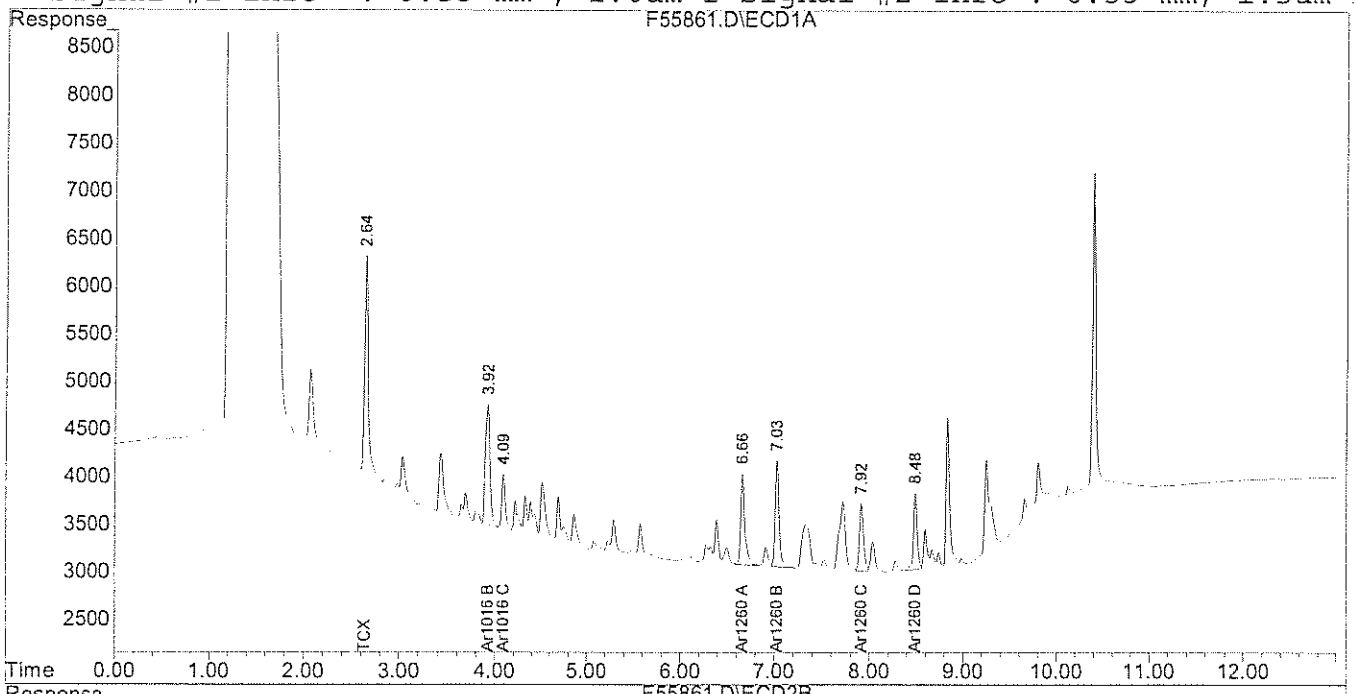
(f)=RT Delta > 1/2 Window (#)=Amounts differ by > 25% (m)=manual int.

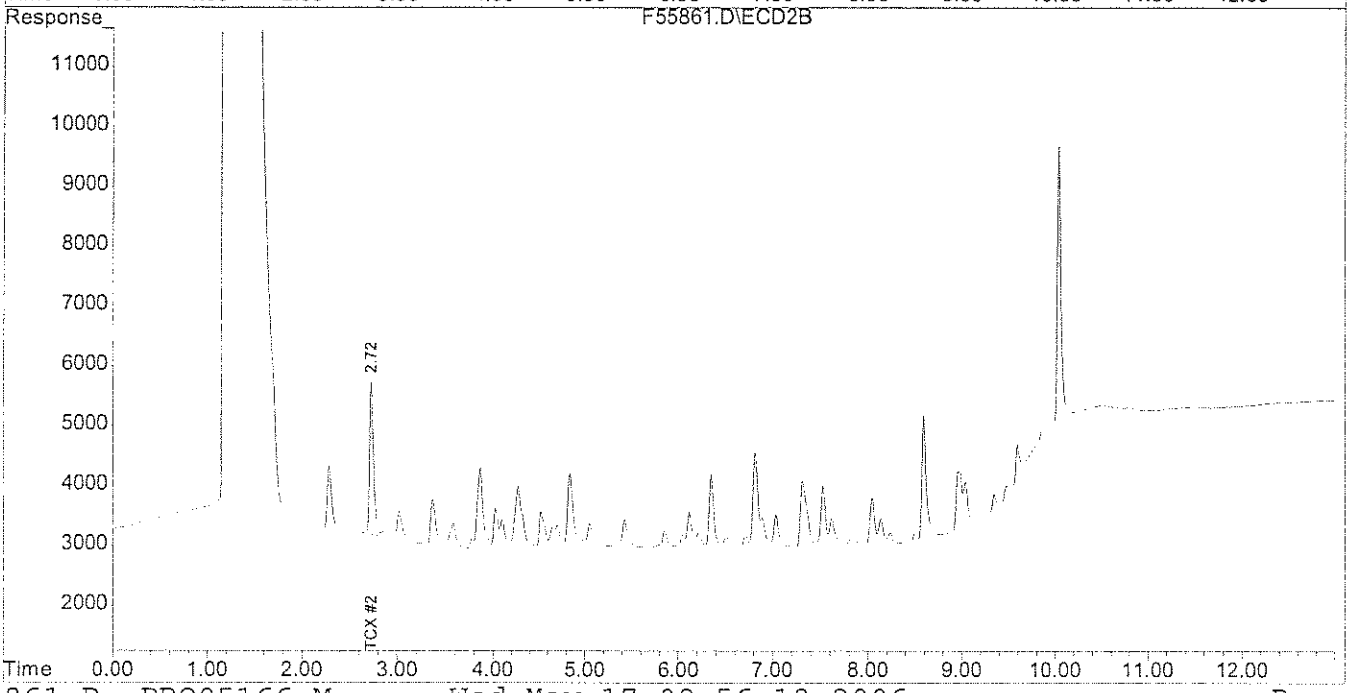
Quantitation Report (QT Reviewed)

Signal #1 : C:\HPCHEM\2\DATA\051606-F\F55861.D\ECD1A.CH Vial: 13
Signal #2 : C:\HPCHEM\2\DATA\051606-F\F55861.D\ECD2B.CH
Acq On : 16 May 2006 7:38 pm Operator:
Sample : P4016 16/60 0.01 Inst : Instrumen
Misc : Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 17 9:56 2006 Quant Results File: PBQ05166.RES

Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 09:54:21 2006
Response via : Multiple Level Calibration
DataAcq Meth : PSTQUICK.M

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film



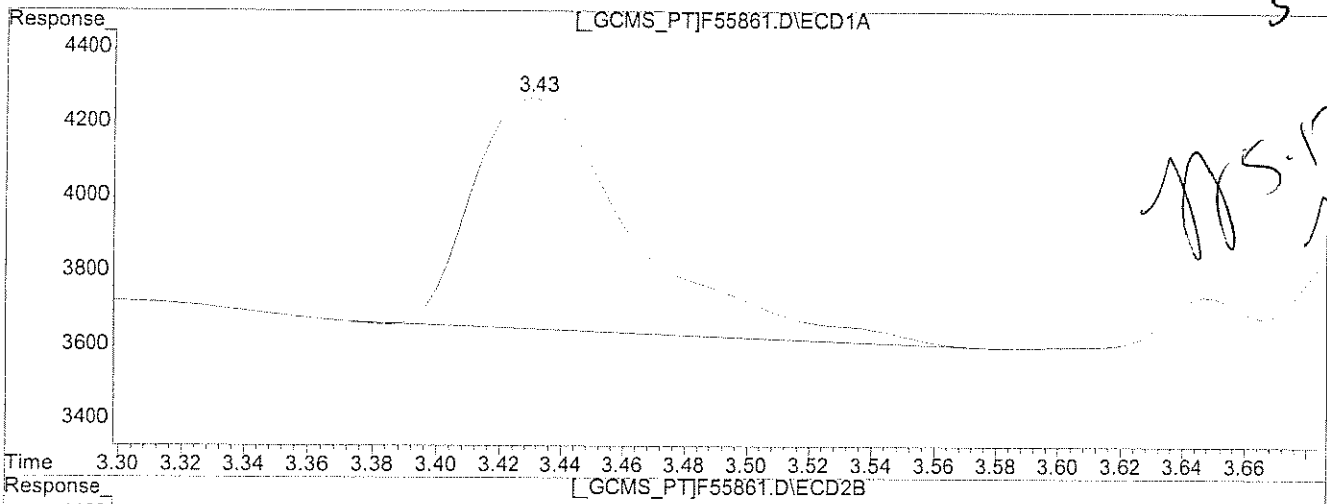


Quantitation Report (Qedit)

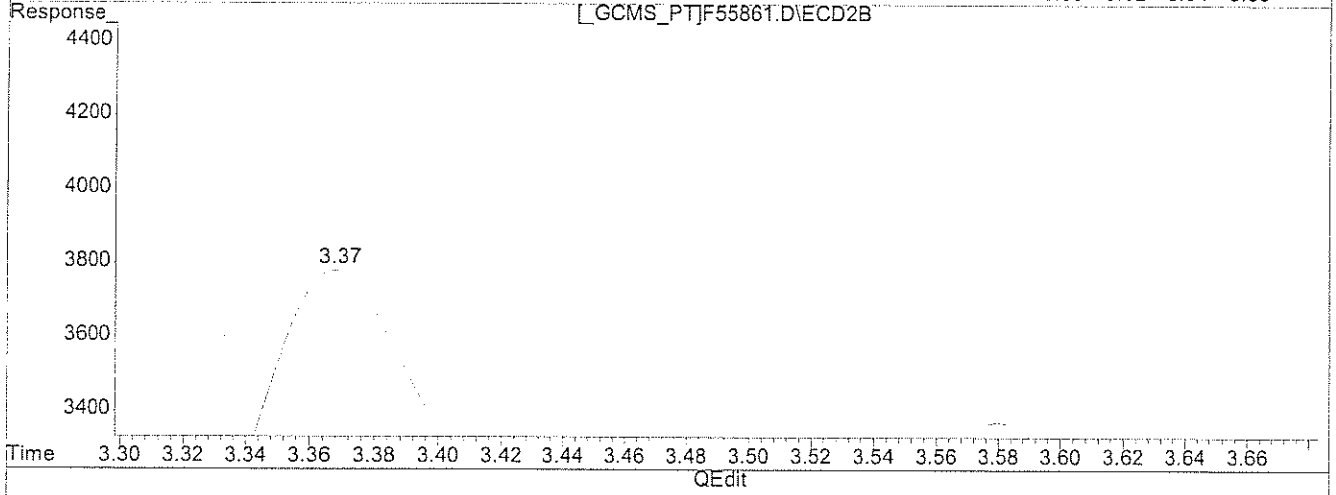
Signal #1 : C:\HPCHEM\2\DATA\051606-F\F55861.D\ECD1A.CH Vial: 13
Signal #2 : C:\HPCHEM\2\DATA\051606-F\F55861.D\ECD2B.CH
Acq On : 16 May 2006 7:38 pm Operator:
Sample : P4016 16/60 0.01 Inst : Instrumen
Misc : Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 17 9:54 2006 Quant Results File: PBQ05166.RES

Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 09:54:21 2006
Response via : Multiple Level Calibration

5.17.06



5.17.06 Shoulder



(2) Ar1016 A (L1)

R.T.	Response	Conc
3.43	22374	-0.00
3.92	43320	0.00
4.09	15699	0.00

(2) Ar1016 A #2 (L1)

R.T.	Response	Conc
3.37	27177	-0.01
3.86	57040	-0.01
4.03	21986	-0.01

(+) = Expected Retention Time

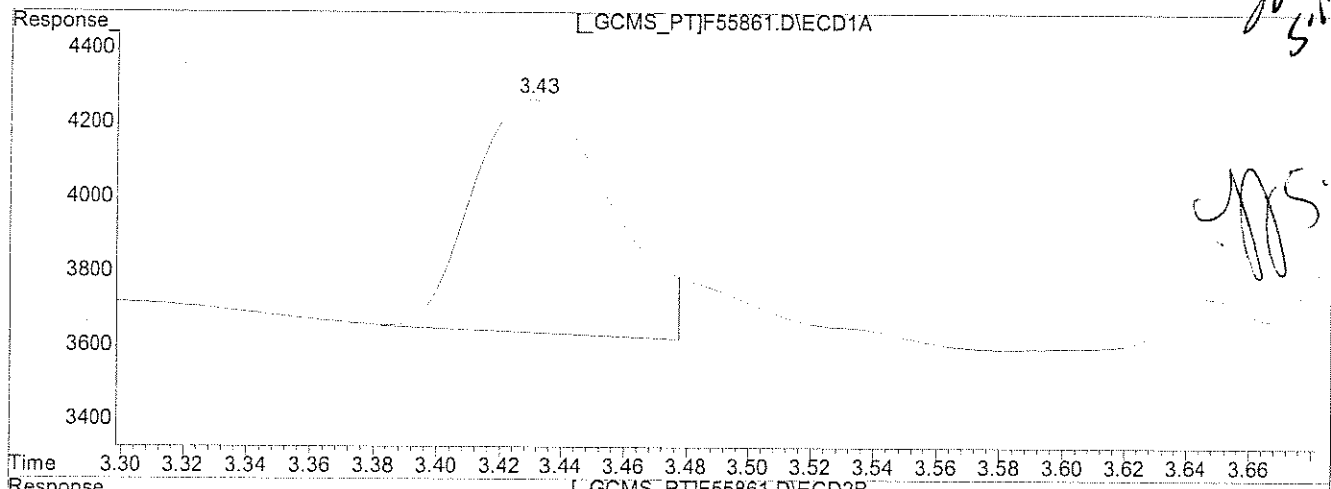
Quantitation Report (Qedit)

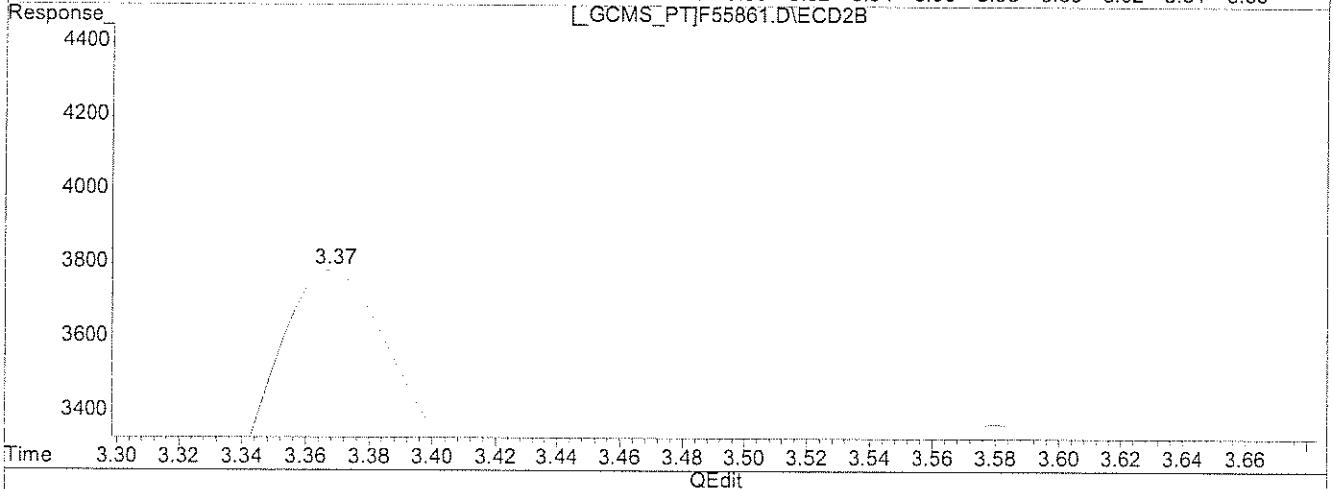
Signal #1 : C:\HPCHEM\2\DATA\051606-F\F55861.D\ECD1A.CH Vial: 13
Signal #2 : C:\HPCHEM\2\DATA\051606-F\F55861.D\ECD2B.CH
Acq On : 16 May 2006 7:38 pm Operator:
Sample : P4016 16/60 0.01 Inst : Instrumen
Misc : Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 17 9:54 2006 Quant Results File: PBQ05166.RES

Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 09:54:21 2006
Response via : Multiple Level Calibration

5.17.06

5.17.06





(2) Ar1016 A #2 (L1)

R.T.	Response	Conc
3.43	19531	-0.00
3.92	43320	0.00
4.09	15699	0.00

(2) Ar1016 A #2 (L1)

R.T.	Response	Conc
3.37	27177	-0.01
3.86	57040	-0.01
4.03	21986	-0.01

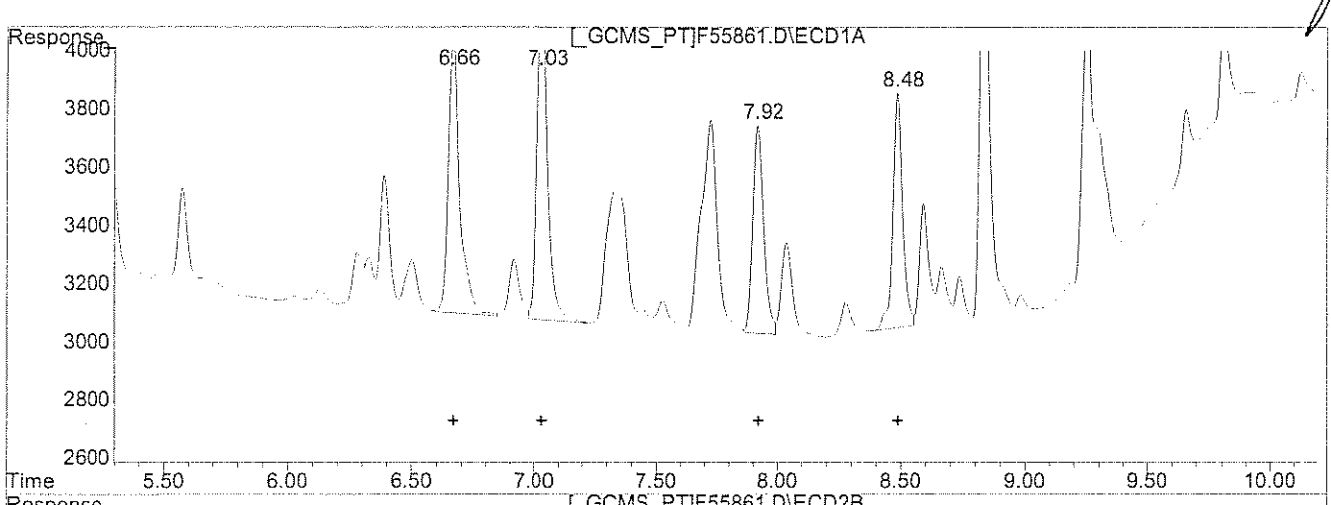
(+) = Expected Retention Time

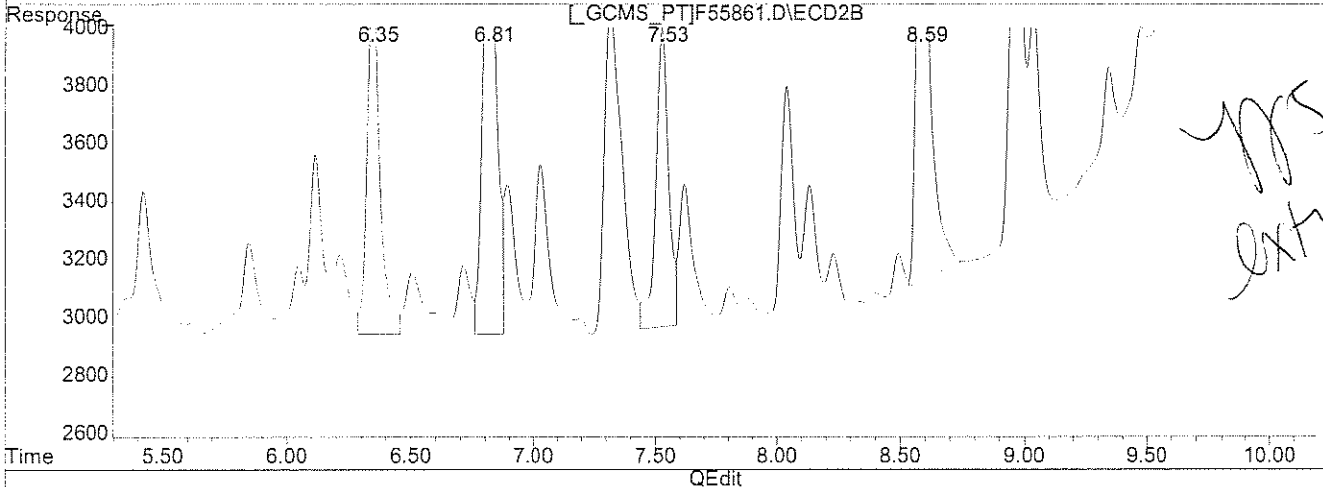
Quantitation Report (Qedit)

Signal #1 : C:\HPCHEM\2\DATA\051606-F\F55861.D\ECD1A.CH Vial: 13
Signal #2 : C:\HPCHEM\2\DATA\051606-F\F55861.D\ECD2B.CH
Acq On : 16 May 2006 7:38 pm Operator:
Sample : P4016 16/60 0.01 Inst : Instrumen
Misc : Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 17 9:54 2006 Quant Results File: PBQ05166.RES

Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 09:54:21 2006
Response via : Multiple Level Calibration

5-17-06





5-17-06 extra ana

(5) Ar1260 A (L2)

R.T.	Response	Conc
6.67	30275	0.00
7.03	33089	0.00
7.92	22273	0.01
8.49	22028	0.00

(5) Ar1260 A #2 (L2)

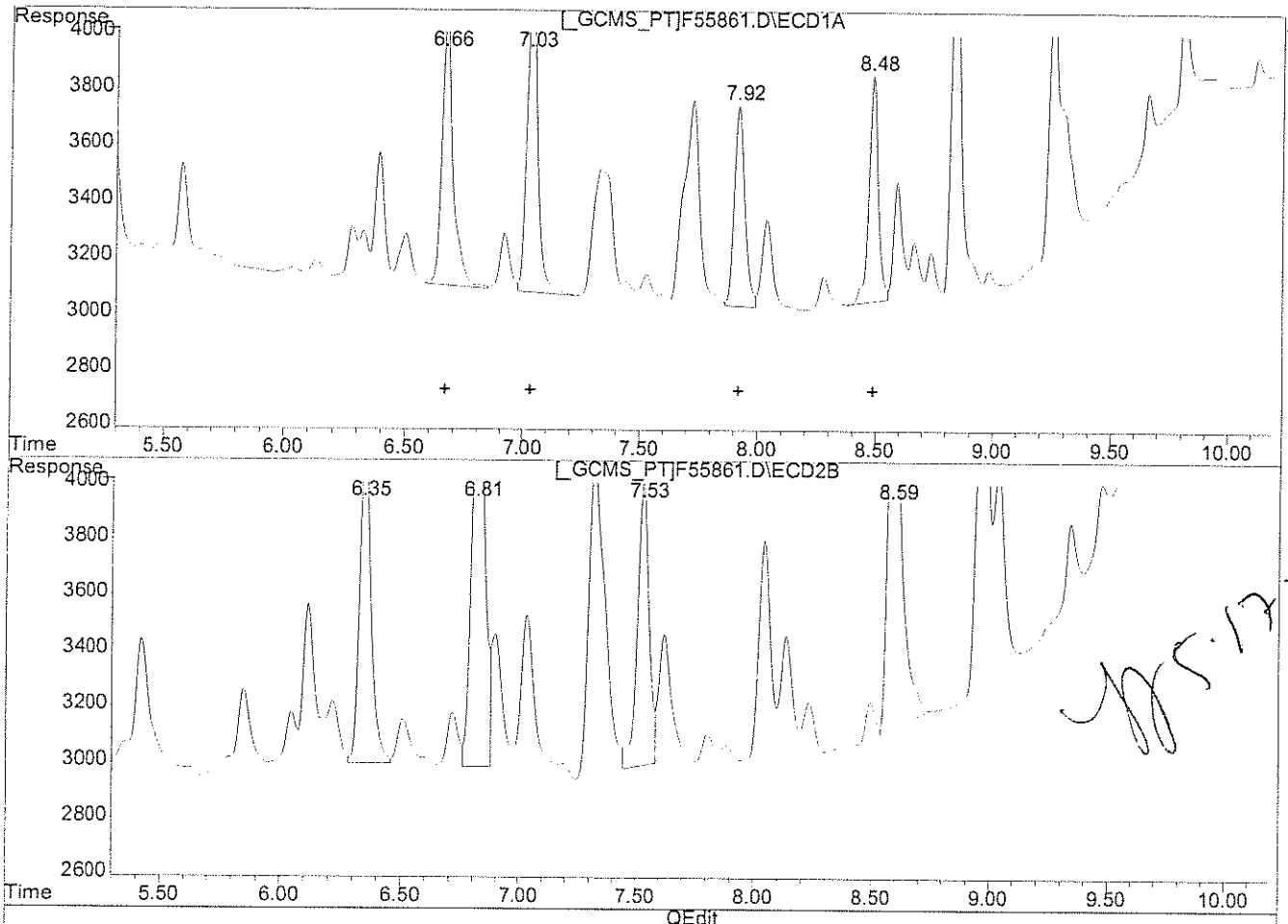
R.T.	Response	Conc
6.35	45036	-0.02
6.82	61803	-0.01
7.53	37673	-0.01
8.59	68293	-0.00

(+) = Expected Retention Time

Quantitation Report (Qedit)

Signal #1 : C:\HPCHEM\2\DATA\051606-F\F55861.D\ECD1A.CH Vial: 13
Signal #2 : C:\HPCHEM\2\DATA\051606-F\F55861.D\ECD2B.CH
Acq On : 16 May 2006 7:38 pm Operator:
Sample : P4016 16/60 0.01 Inst : Instrumen
Misc : Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 17 9:54 2006 Quant Results File: PBQ05166.RES

Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 09:54:21 2006
Response via : Multiple Level Calibration



(5) Ar1260 A #2 (L2)

R.T.	Response	Conc
6.67	30275	0.00
7.03	33089	0.00
7.92	22273	0.01
8.49	22028	0.00

(5) Ar1260 A #2 (L2)

R.T.	Response	Conc
6.35	39717	-0.02
6.81	60122	-0.01
7.53	35850	-0.01
8.59	68293	-0.00

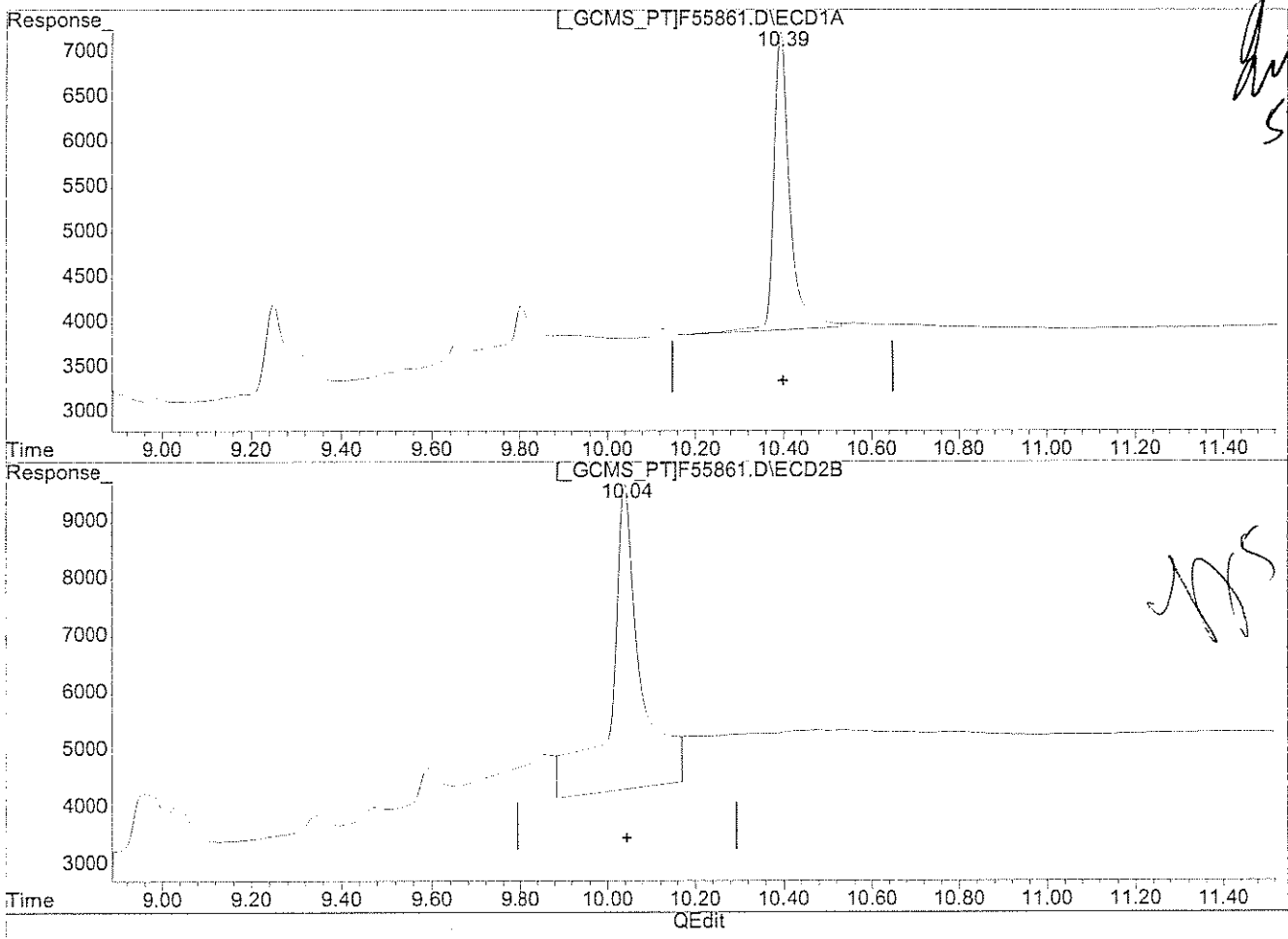
(+) = Expected Retention Time

F55861.D PBQ05166.M Wed May 17 09:55:59 2006

Quantitation Report (Qedit)

Signal #1 : C:\HPCHEM\2\DATA\051606-F\F55861.D\ECD1A.CH Vial: 13
Signal #2 : C:\HPCHEM\2\DATA\051606-F\F55861.D\ECD2B.CH
Acq On : 16 May 2006 7:38 pm Operator:
Sample : P4016 16/60 0.01 Inst : Instrumen
Misc : Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 17 9:54 2006 Quant Results File: PBQ05166.RES

Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 09:54:21 2006
Response via : Multiple Level Calibration



5.17.06

5.17.06 baseline

(9) DCB (S)

10.39min -0.001ug/ml

response 81717

(9) DCB #2 (S)

10.04min -0.000ug/ml

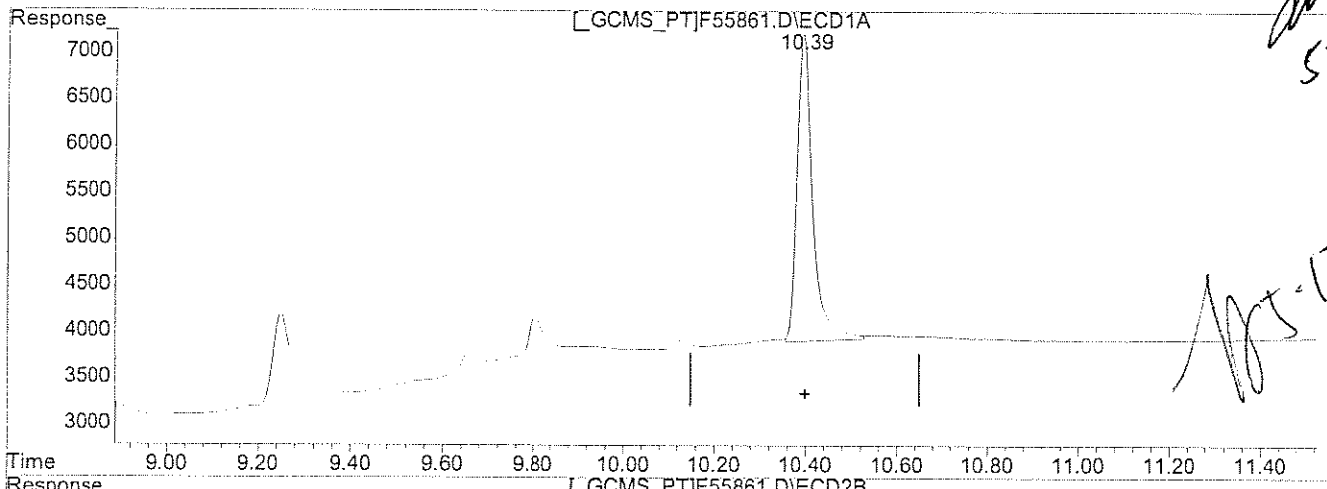
response 257756

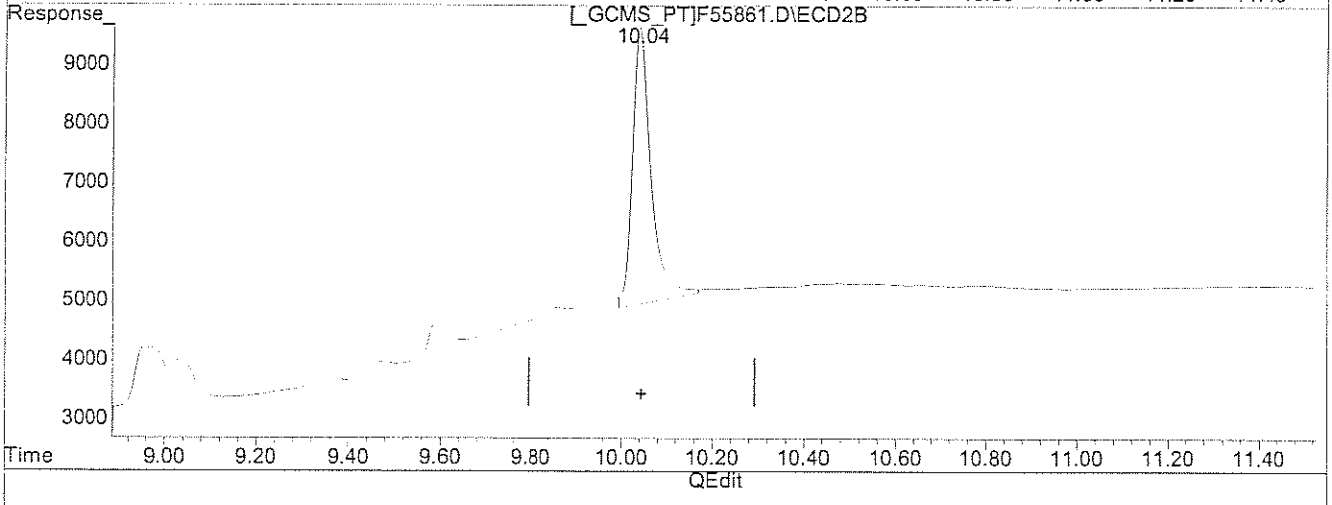
(+) = Expected Retention Time

Quantitation Report (Qedit)

Signal #1 : C:\HPCHEM\2\DATA\051606-F\F55861.D\ECD1A.CH Vial: 13
Signal #2 : C:\HPCHEM\2\DATA\051606-F\F55861.D\ECD2B.CH
Acq On : 16 May 2006 7:38 pm Operator:
Sample : P4016 16/60 0.01 Inst : Instrumen
Misc : Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 17 9:54 2006 Quant Results File: PBQ05166.RES

Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 09:54:21 2006
Response via : Multiple Level Calibration





QEdit

(9) DCB (S)

10.39min -0.001ug/ml m

response 79033

(9) DCB #2 (S)

10.04min -0.003ug/ml m

response 133576

(+) = Expected Retention Time

F55861.D PBQ05166.M Wed May 17 09:56:13 2006

Quantitation Report (QT Reviewed)

Signal #1 : C:\HPCHEM\2\DATA\051606-F\F55862.D\ECD1A.CH Vial: 14
Signal #2 : C:\HPCHEM\2\DATA\051606-F\F55862.D\ECD2B.CH
Acq On : 16 May 2006 7:53 pm Operator:
Sample : P4017 16/60 0.02 Inst : Instrumen
Misc : Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 17 9:56 2006 Quant Results File: PBQ05166.RES

Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 09:54:21 2006
Response via : Initial Calibration
DataAcq Meth : PSTQUICK.M

5-17-06 5-17-06

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film

Compound	RT#1	RT#2	Resp#1	Resp#2	ug/ml	ug/ml
System Monitoring Compounds						
1) S TCX	2.64	2.73	122756	153548	0.004	0.003
Spiked Amount 0.100	Range 46 - 122			Recovery =	4.00%#	3.00%#
9) S DCB	10.40	10.04	155614	248581	0.001m	N.D. m#
Spiked Amount 0.100	Range 26 - 86			Recovery =	1.00%#	0.00%#
Target Compounds						
2) L1 Ar1016 A	3.44	3.37	42784	51825	0.011	0.006 #
3) L1 Ar1016 B	3.93	3.87	80946	111881	0.014	0.008 #
4) L1 Ar1016 C	4.09	4.03	29570	41472	0.016	0.008 #
Sum Ar1016 A			153300	205178	0.041	0.022
Average Ar1016 A					0.014	0.007
5) L2 Ar1260 A	6.67	6.35	53627	81296	0.014	0.002 #
6) L2 Ar1260 B	7.04	6.82	59571	113223	0.013	0.004 #
7) L2 Ar1260 C	7.92	7.54	42912	71915	0.017	0.007 #
8) L2 Ar1260 D	8.49	8.59	42372	139041	0.016	0.012 #
Sum Ar1260 A			198483	405475	0.060	0.024
Average Ar1260 A					0.015	0.006

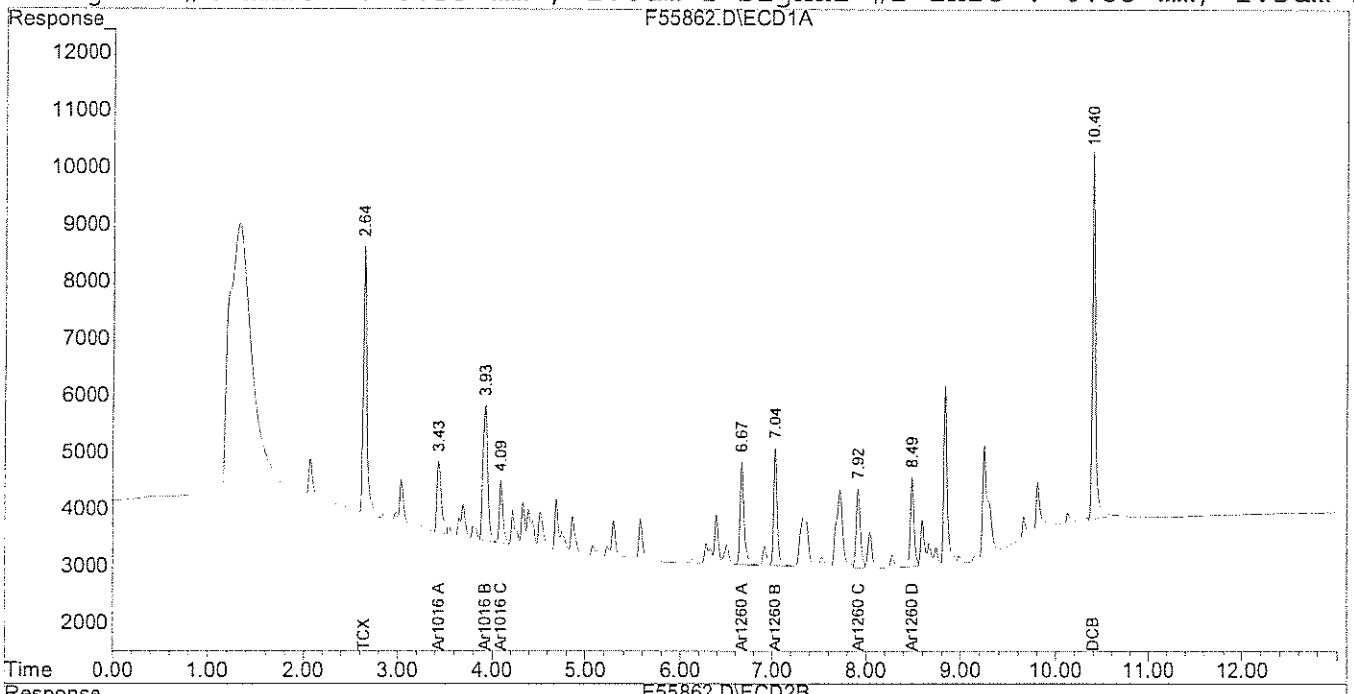
(f)=RT Delta > 1/2 Window (#)=Amounts differ by > 25% (m)=manual int.

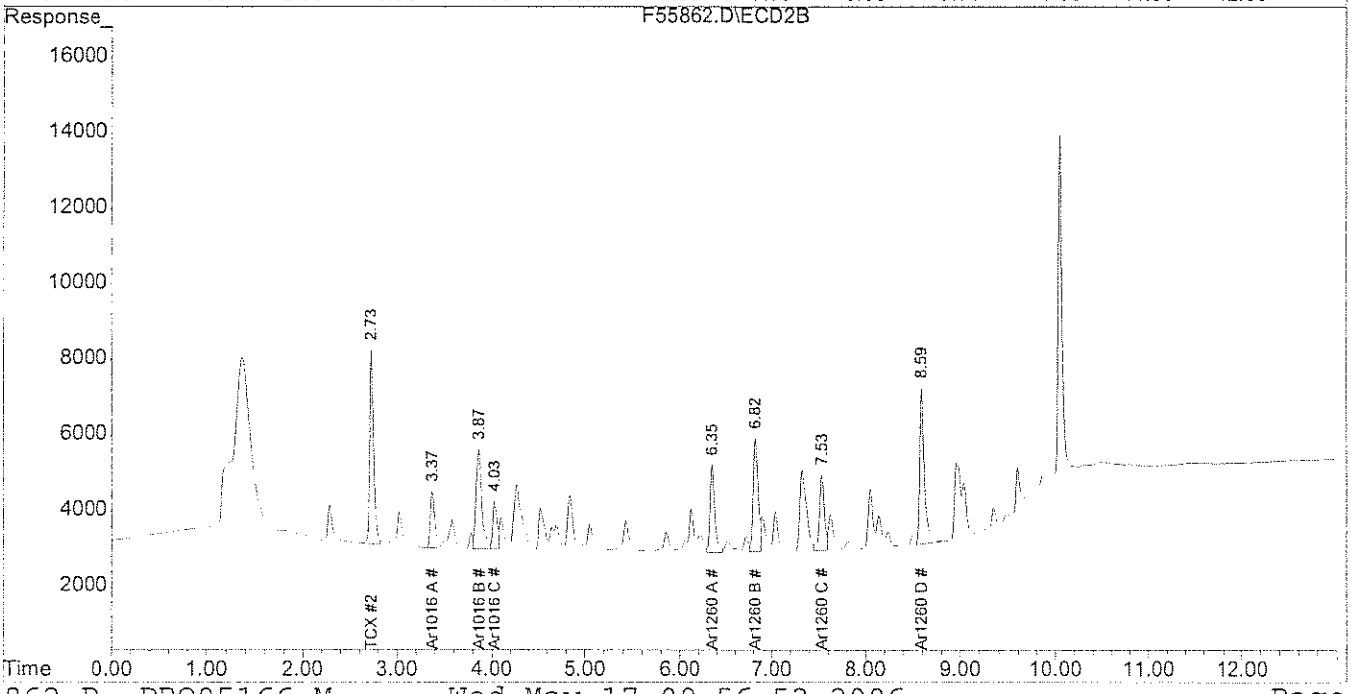
Quantitation Report (QT Reviewed)

Signal #1 : C:\HPCHEM\2\DATA\051606-F\F55862.D\ECD1A.CH Vial: 14
Signal #2 : C:\HPCHEM\2\DATA\051606-F\F55862.D\ECD2B.CH
Acq On : 16 May 2006 7:53 pm Operator:
Sample : P4017 16/60 0.02 Inst : Instrumen
Misc : Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 17 9:56 2006 Quant Results File: PBQ05166.RES

Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 09:54:21 2006
Response via : Multiple Level Calibration
DataAcq Meth : PSTQUICK.M

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film

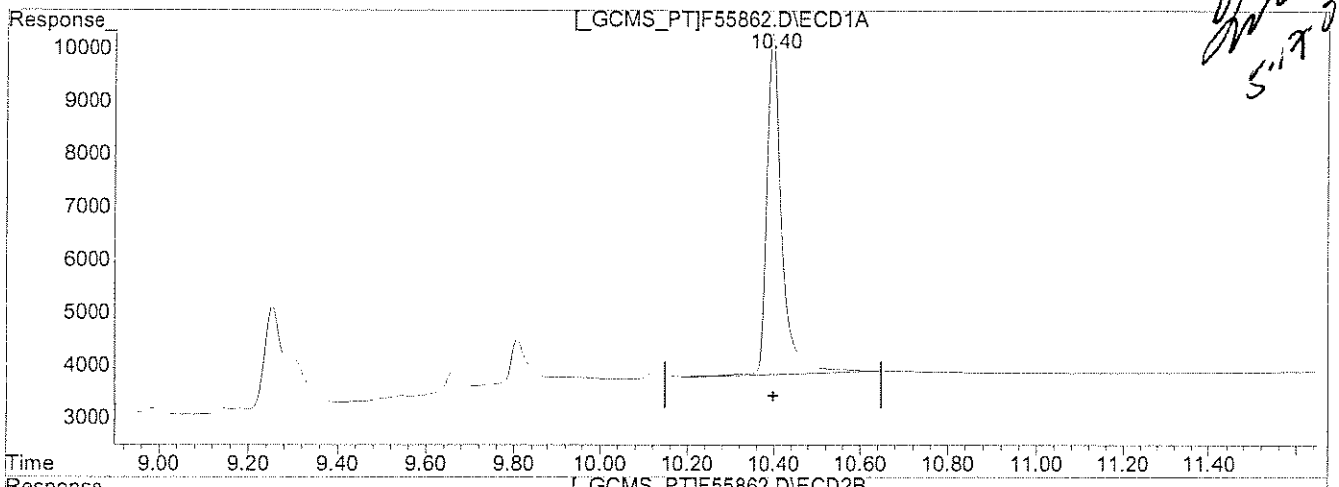


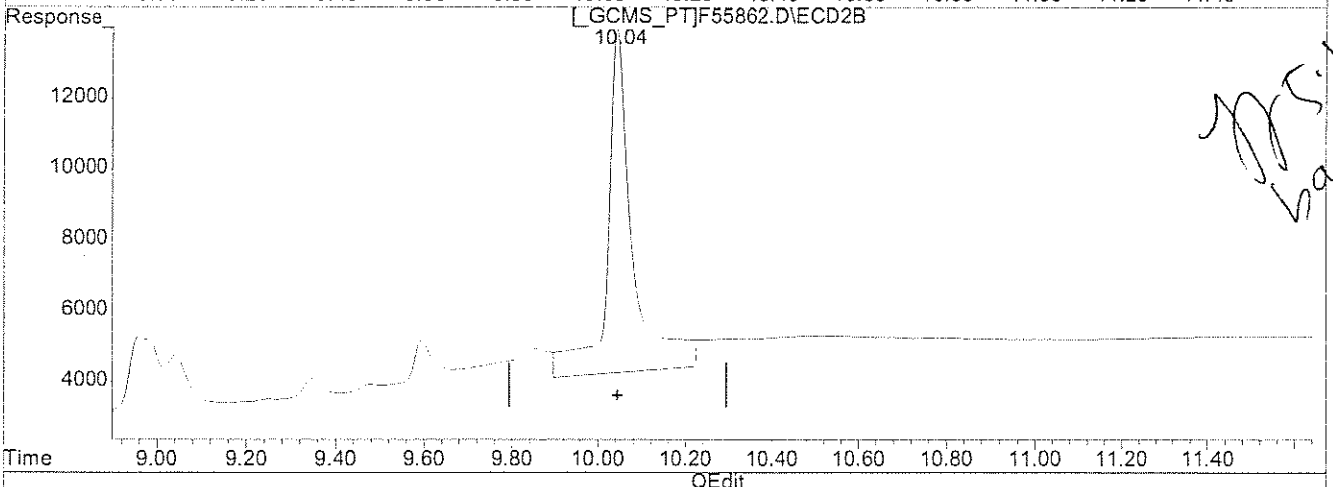


Quantitation Report (Qedit)

Signal #1 : C:\HPCHEM\2\DATA\051606-F\F55862.D\ECD1A.CH Vial: 14
Signal #2 : C:\HPCHEM\2\DATA\051606-F\F55862.D\ECD2B.CH
Acq On : 16 May 2006 7:53 pm Operator:
Sample : P4017 16/60 0.02 Inst : Instrumen
Misc : Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 17 9:56 2006 Quant Results File: PBQ05166.RES

Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 09:54:21 2006
Response via : Multiple Level Calibration





QEdit

(9) DCB (S)

10.40min 0.002ug/ml

response 156938

(9) DCB #2 (S)

10.04min 0.003ug/ml

response 389832

(+) = Expected Retention Time

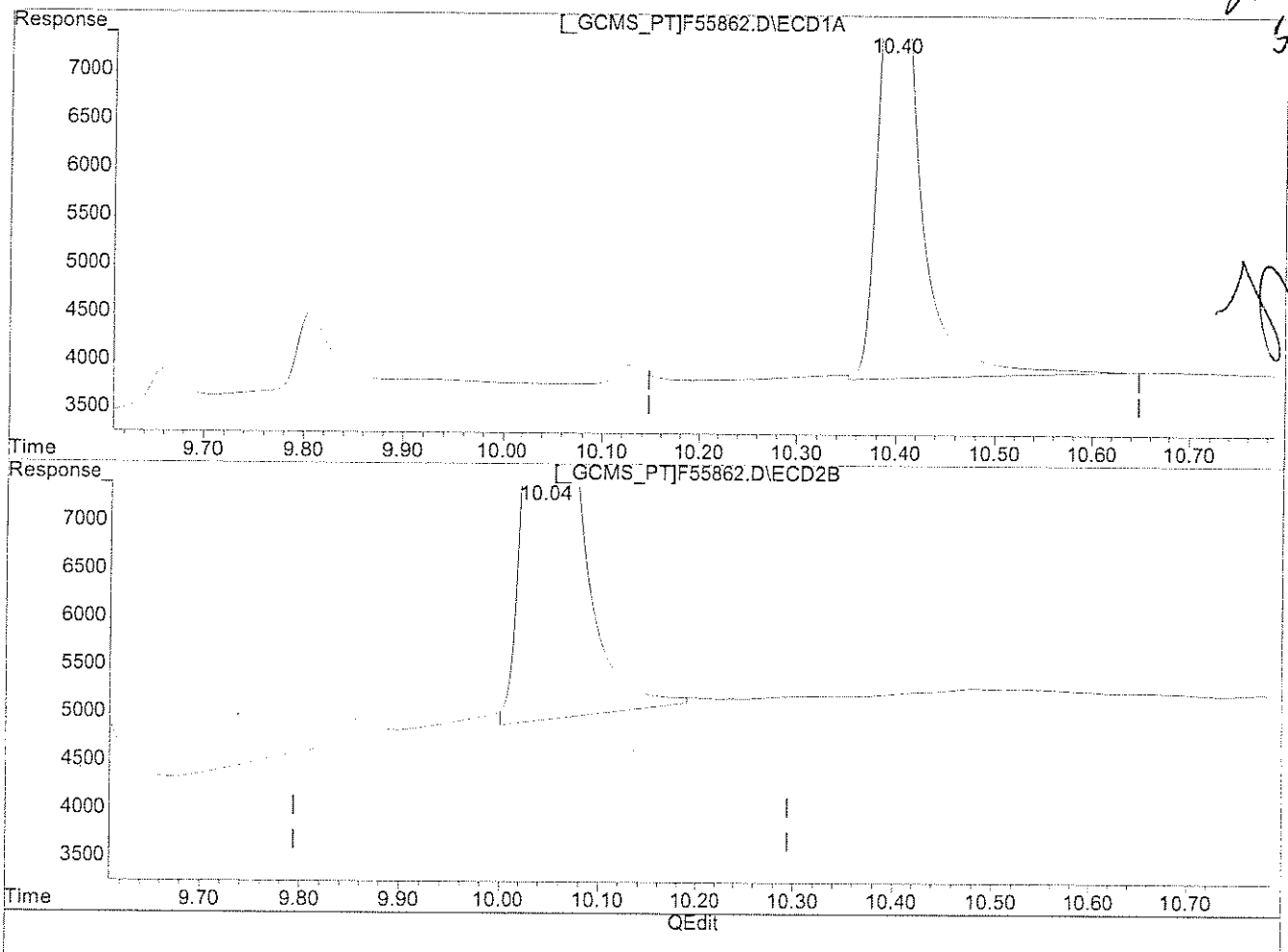
F55862.D PBO05166.M Wed May 17 09:56:36 2006

Quantitation Report (Qedit)

Signal #1 : C:\HPCHEM\2\DATA\051606-F\F55862.D\ECD1A.CH Vial: 14
Signal #2 : C:\HPCHEM\2\DATA\051606-F\F55862.D\ECD2B.CH
Acq On : 16 May 2006 7:53 pm Operator:
Sample : P4017 16/60 0.02 Inst : Instrumen
Misc : Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 17 9:56 2006 Quant Results File: PBQ05166.RES

Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 09:54:21 2006
Response via : Multiple Level Calibration

5.17.06



5.17.06

(9) DCB (S)

10.40min 0.001ug/ml m

response 155614

(9) DCB #2 (S)

10.04min -0.000ug/ml m

response 248581

(+) = Expected Retention Time

Quantitation Report (QT Reviewed)

Signal #1 : C:\HPCHEM\2\DATA\051606-F\F55863.D\ECD1A.CH Vial: 15
Signal #2 : C:\HPCHEM\2\DATA\051606-F\F55863.D\ECD2B.CH
Acq On : 16 May 2006 8:07 pm Operator:
Sample : P4018 16/60 0.05 Inst : Instrumen
Misc : Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 17 10:04 2006 Quant Results File: PBQ05166.RES

Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 09:54:21 2006
Response via : Initial Calibration
DataAcq Meth : PSTQUICK.M

5.17.06 5.17.06

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film

Compound	RT#1	RT#2	Resp#1	Resp#2	ug/ml	ug/ml
System Monitoring Compounds						
1) S TCX	2.64	2.73	328762	401821	0.010	0.010
Spiked Amount 0.100	Range 46 - 122		Recovery =		10.00%#	10.00%#
9) S DCB	10.40	10.05	369688	589584	0.009m	0.009m
Spiked Amount 0.100	Range 26 - 86		Recovery =		9.00%#	9.00%#
Target Compounds						
2) L1 Ar1016 A	3.44	3.37	102888	121190	0.051	0.050
3) L1 Ar1016 B	3.93	3.87	187238	263561	0.051	0.050
4) L1 Ar1016 C	4.10	4.03	75426	99033	0.054	0.051
Sum Ar1016 A			365551	483784	0.156	0.151
Average Ar1016 A					0.052	0.050
5) L2 Ar1260 A	6.67	6.35	126226	190696	0.053	0.054
6) L2 Ar1260 B	7.04	6.82	148952	269972	0.053	0.054
7) L2 Ar1260 C	7.92	7.53	105396	183351	0.052	0.056
8) L2 Ar1260 D	8.49	8.59	102899	351344	0.051	0.051
Sum Ar1260 A			483473	995362	0.209	0.216
Average Ar1260 A					0.052	0.054

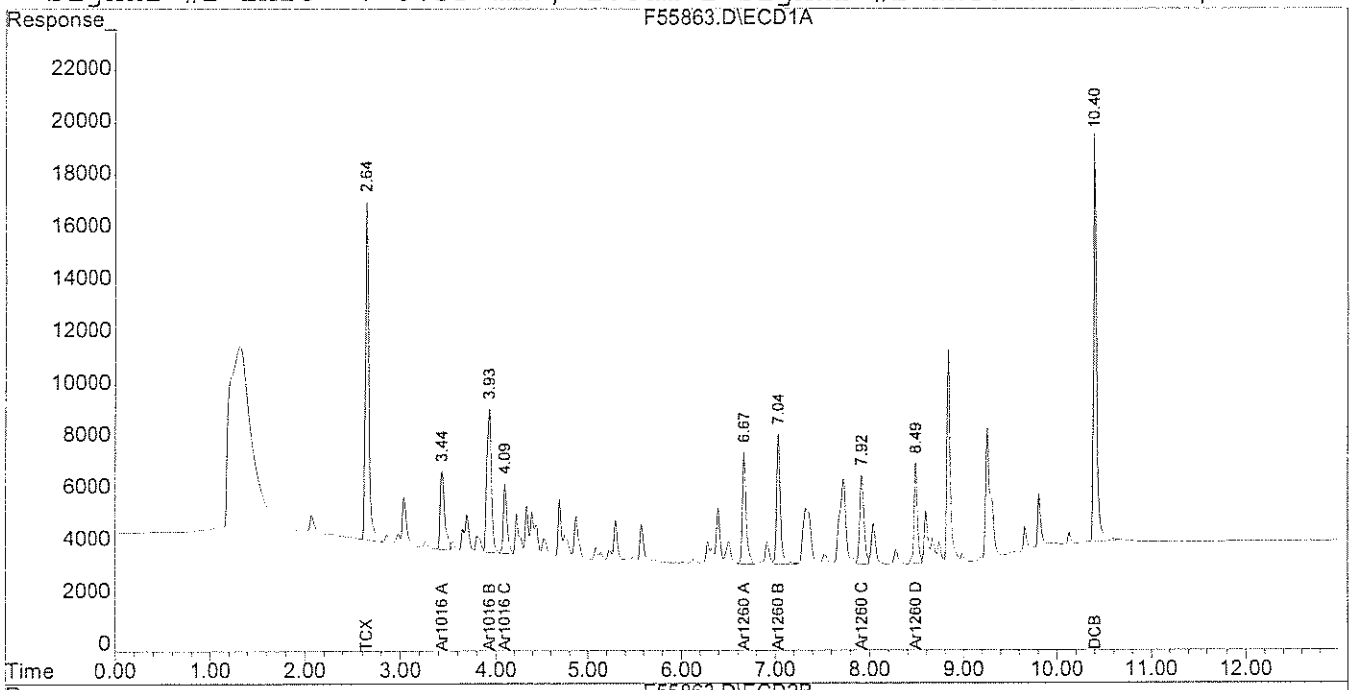
(f)=RT Delta > 1/2 Window (#)=Amounts differ by > 25% (m)=manual int.

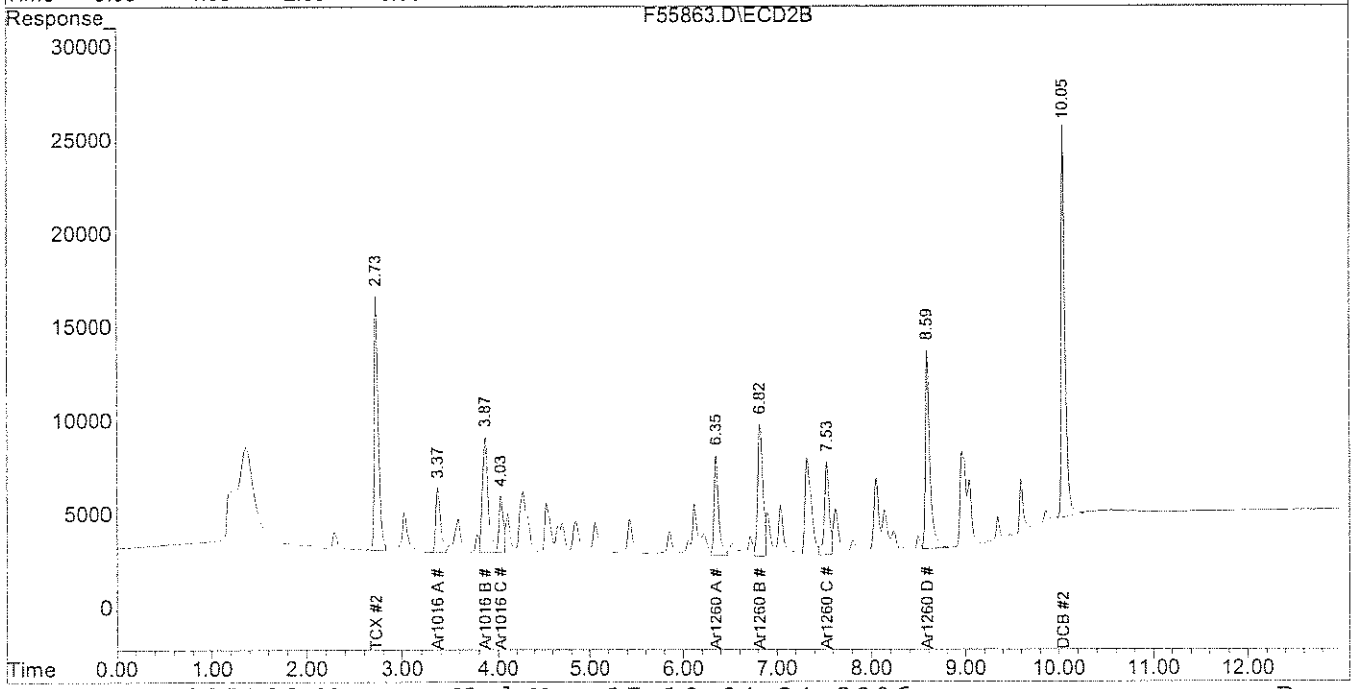
Quantitation Report (QT Reviewed)

Signal #1 : C:\HPCHEM\2\DATA\051606-F\F55863.D\ECD1A.CH Vial: 15
Signal #2 : C:\HPCHEM\2\DATA\051606-F\F55863.D\ECD2B.CH
Acq On : 16 May 2006 8:07 pm Operator:
Sample : P4018 16/60 0.05 Inst : Instrumen
Misc : Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 17 10:04 2006 Quant Results File: PBQ05166.RES

Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 09:54:21 2006
Response via : Multiple Level Calibration
DataAcq Meth : PSTQUICK.M

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film





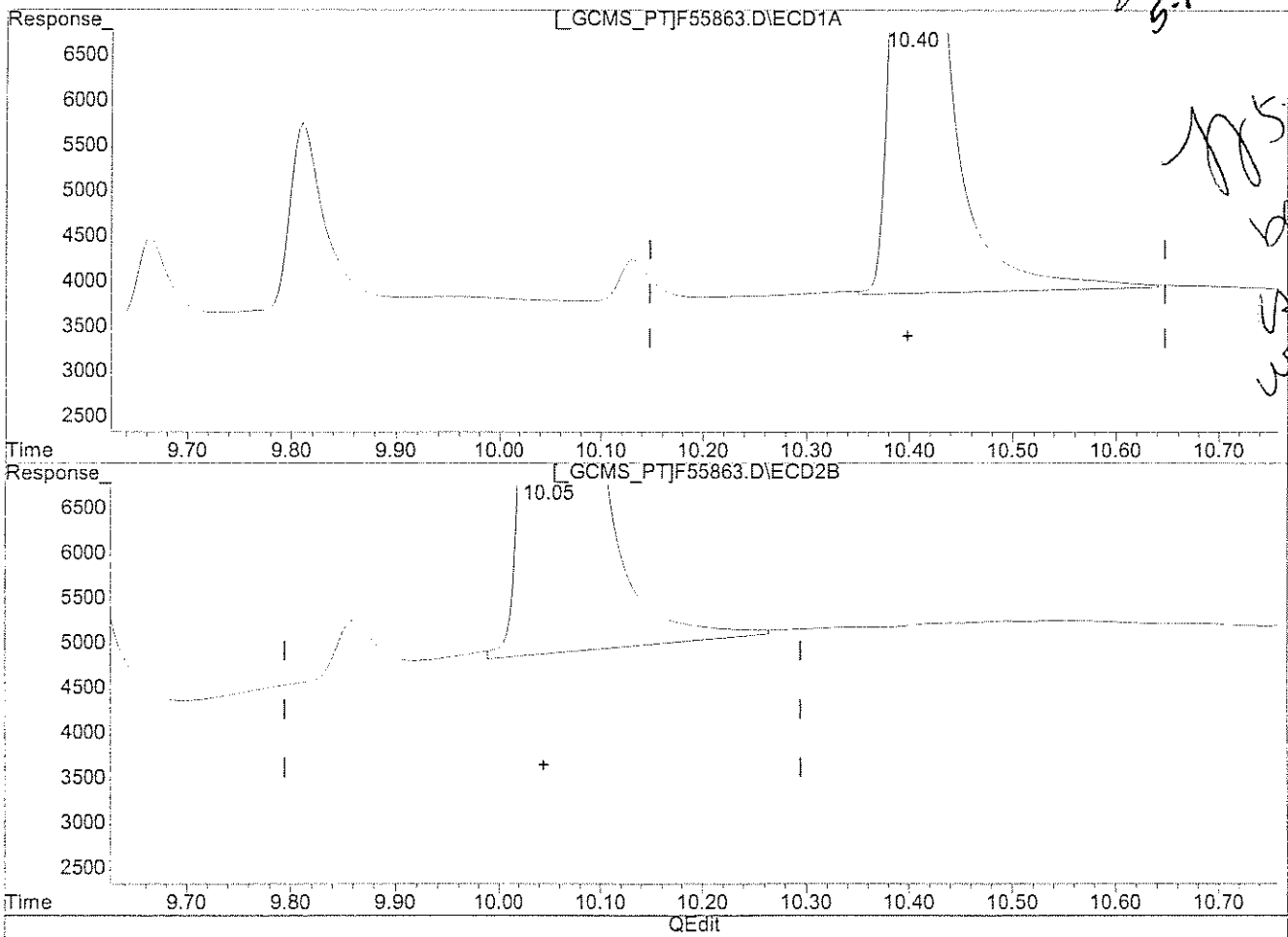
Quantitation Report (Qedit)

Signal #1 : C:\HPCHEM\2\DATA\051606-F\F55863.D\ECD1A.CH Vial: 15
Signal #2 : C:\HPCHEM\2\DATA\051606-F\F55863.D\ECD2B.CH
Acq On : 16 May 2006 8:07 pm Operator:
Sample : P4018 16/60 0.05 Inst : Instrumen
Misc : Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 17 9:57 2006 Quant Results File: PBQ05166.RES

Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 09:54:21 2006
Response via : Multiple Level Calibration

5-17-06

5-17-06 baseline issues were present before manual



(9) DCB (S)

10.40min 0.009ug/ml m

response 369688

(9) DCB #2 (S)

10.05min 0.009ug/ml m

response 589584

(+) = Expected Retention Time

Quantitation Report (QT Reviewed)

Signal #1 : C:\HPCHEM\2\DATA\051606-F\F55864.D\ECD1A.CH Vial: 16
Signal #2 : C:\HPCHEM\2\DATA\051606-F\F55864.D\ECD2B.CH
Acq On : 16 May 2006 8:22 pm Operator:
Sample : P4019 16/60 0.1 Inst : Instrumen
Misc : Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 17 10:04 2006 Quant Results File: PBQ05166.RES

Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 09:54:21 2006
Response via : Initial Calibration
DataAcq Meth : PSTQUICK.M

5-17-06 5-17-06

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film

Compound	RT#1	RT#2	Resp#1	Resp#2	ug/ml	ug/ml
System Monitoring Compounds						
1) S TCX	2.64	2.73	688147	815855	0.022	0.022
Spiked Amount 0.100	Range 46 - 122		Recovery =		22.00%#	22.00%#
9) S DCB	10.40	10.05	698690	1099353	0.020	0.022m
Spiked Amount 0.100	Range 26 - 86		Recovery =		20.00%#	22.00%#
Target Compounds						
2) L1 Ar1016 A	3.43	3.37	199668	227033	0.116	0.117
3) L1 Ar1016 B	3.93	3.87	361008	502332	0.111	0.117
4) L1 Ar1016 C	4.09	4.03	146603	188678	0.113	0.118
Sum Ar1016 A			707279	918043	0.340	0.352
Average Ar1016 A					0.113	0.117
5) L2 Ar1260 A	6.67	6.35	234823	325259	0.112	0.119
6) L2 Ar1260 B	7.04	6.82	274639	476109	0.109	0.121
7) L2 Ar1260 C	7.92	7.53	203415	325317	0.107	0.119
8) L2 Ar1260 D	8.49	8.59	199734	691593	0.106	0.114
Sum Ar1260 A			912612	1818277	0.433	0.473
Average Ar1260 A					0.108	0.118

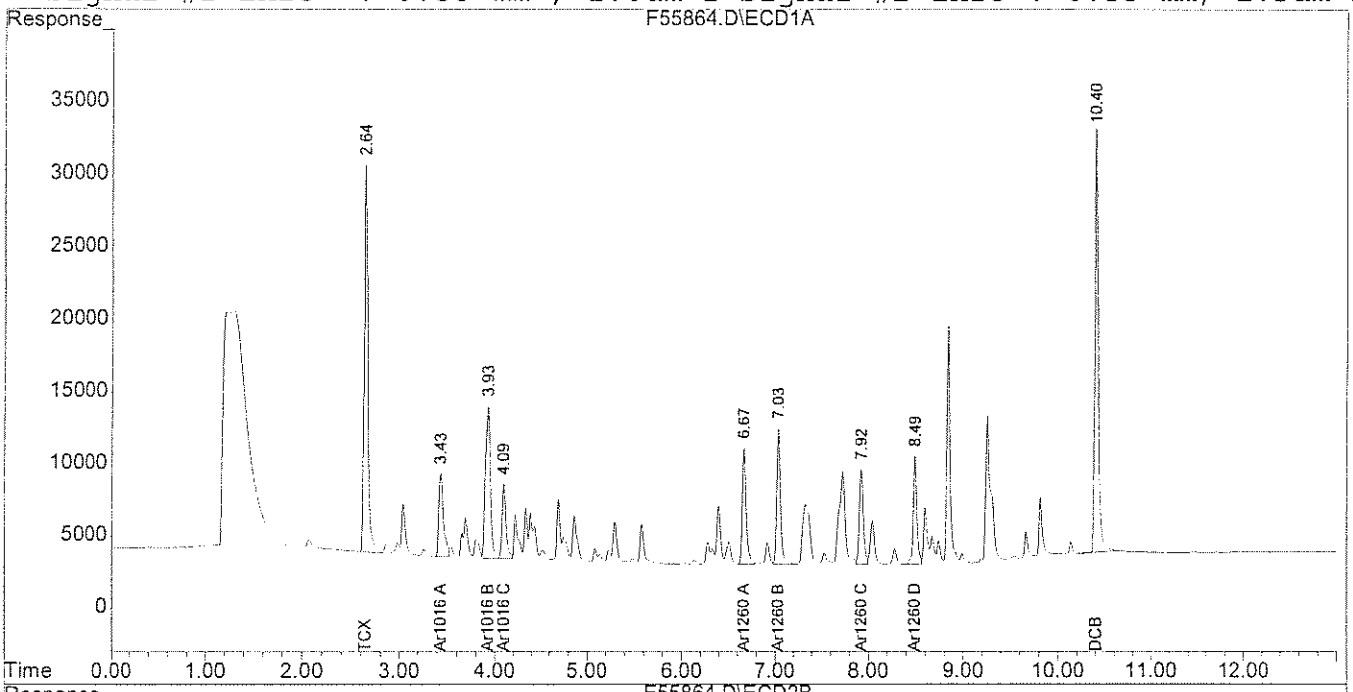
(f)=RT Delta > 1/2 Window (#)=Amounts differ by > 25% (m)=manual int.

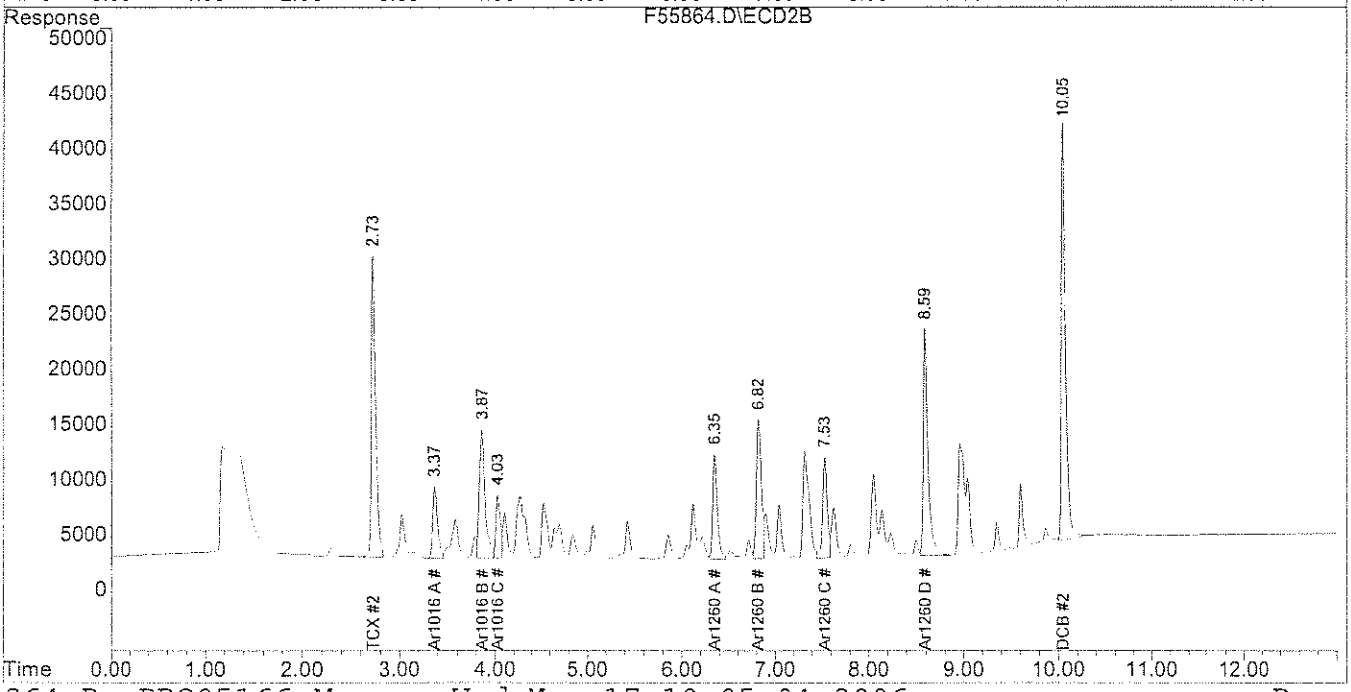
Quantitation Report (QT Reviewed)

Signal #1 : C:\HPCHEM\2\DATA\051606-F\F55864.D\ECD1A.CH Vial: 16
Signal #2 : C:\HPCHEM\2\DATA\051606-F\F55864.D\ECD2B.CH
Acq On : 16 May 2006 8:22 pm Operator:
Sample : P4019 16/60 0.1 Inst : Instrumen
Misc : Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 17 10:04 2006 Quant Results File: PBQ05166.RES

Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 09:54:21 2006
Response via : Multiple Level Calibration
DataAcq Meth : PSTQUICK.M

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film



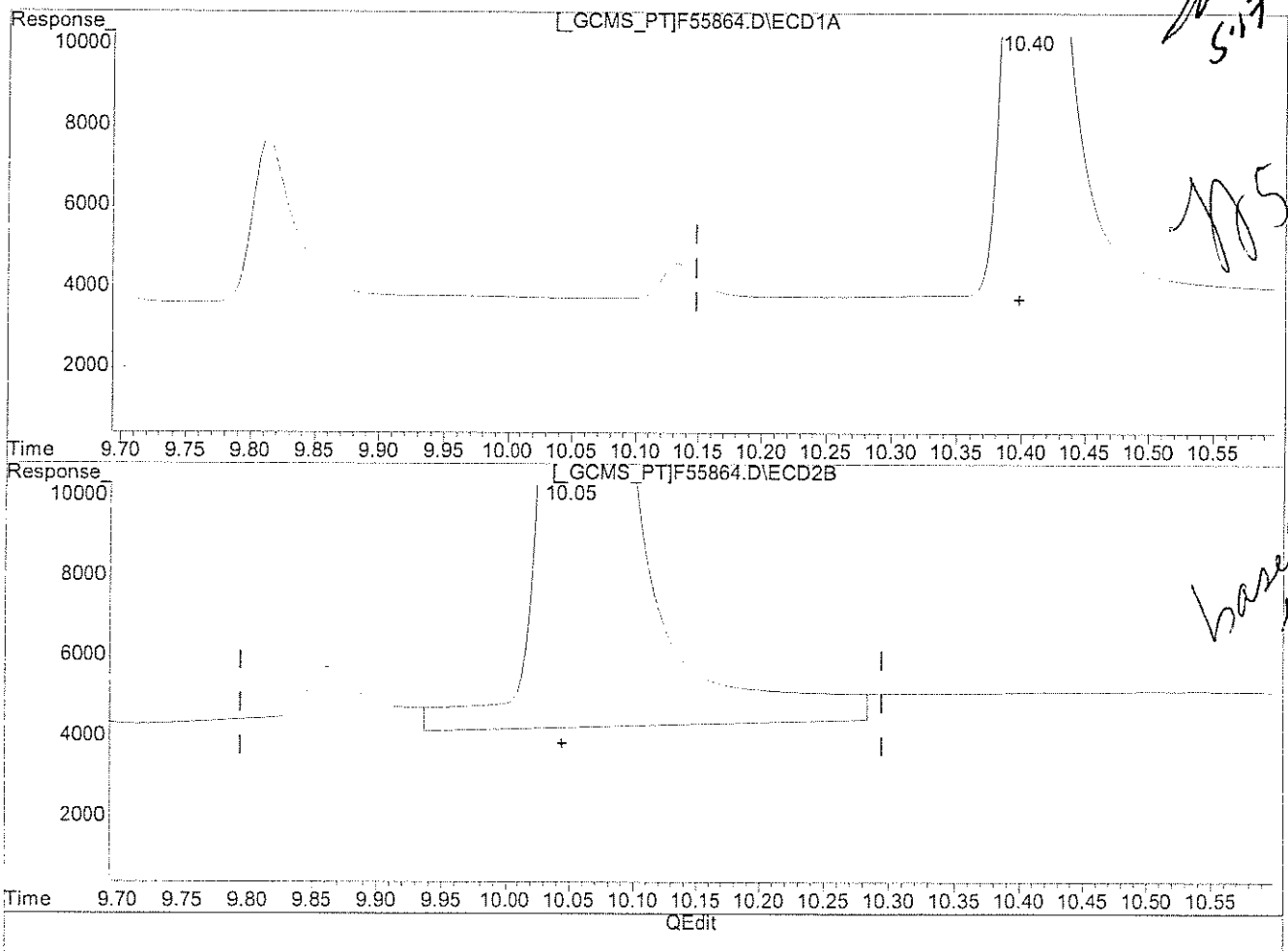


Quantitation Report (Qedit)

Signal #1 : C:\HPCHEM\2\DATA\051606-F\F55864.D\ECD1A.CH Vial: 16
Signal #2 : C:\HPCHEM\2\DATA\051606-F\F55864.D\ECD2B.CH
Acq On : 16 May 2006 8:22 pm Operator:
Sample : P4019 16/60 0.1 Inst : Instrumen
Misc : Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 17 10:04 2006 Quant Results File: PBQ05166.RES

Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 09:54:21 2006
Response via : Multiple Level Calibration

5.17.06



5.17.06

baseline/ extra area

(9) DCB (S)

10.40min 0.020ug/ml

response 698690

(9) DCB #2 (S)

10.05min 0.025ug/ml

response 1203022

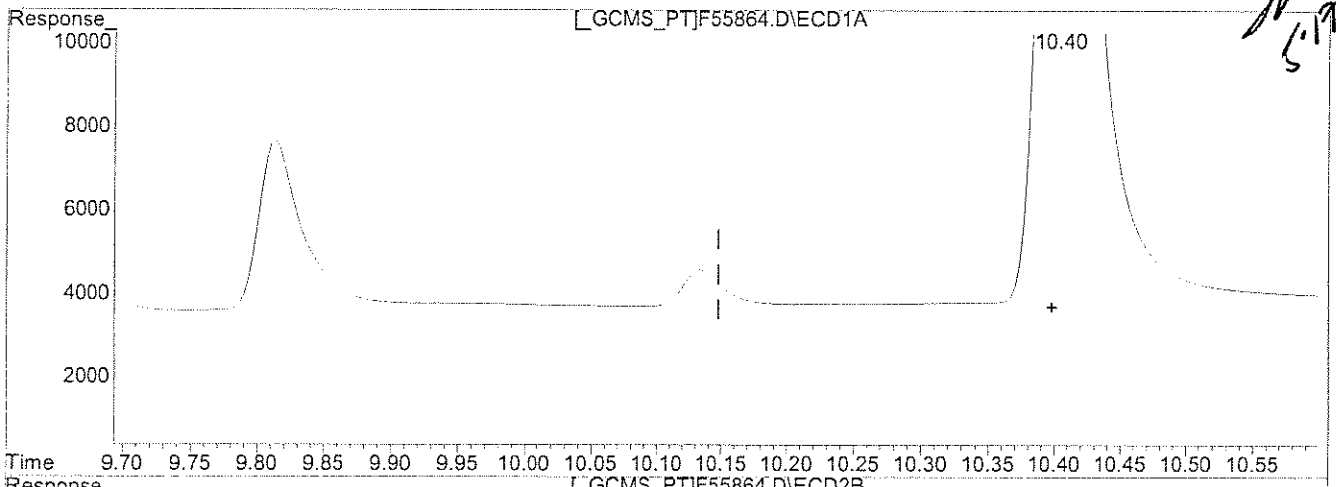
(+) = Expected Retention Time

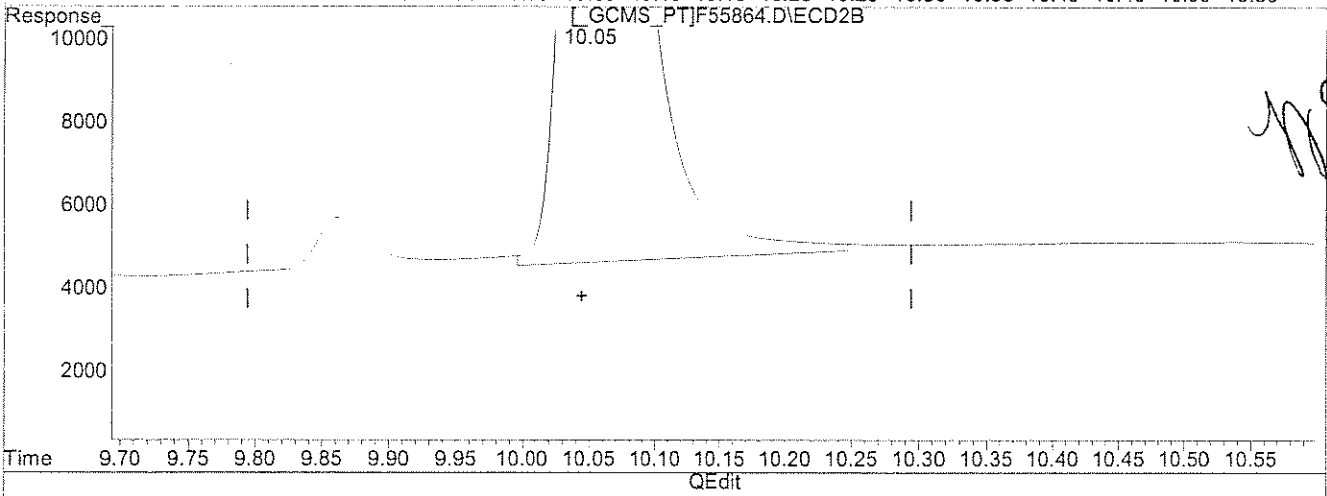
Quantitation Report (Qedit)

Signal #1 : C:\HPCHEM\2\DATA\051606-F\F55864.D\ECD1A.CH Vial: 16
Signal #2 : C:\HPCHEM\2\DATA\051606-F\F55864.D\ECD2B.CH
Acq On : 16 May 2006 8:22 pm Operator:
Sample : P4019 16/60 0.1 Inst : Instrumen
Misc : Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 17 10:04 2006 Quant Results File: PBQ05166.RES

Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 09:54:21 2006
Response via : Multiple Level Calibration

5.17.06





5.17.06

(9) DCB (S)

10.40min 0.020ug/ml

response 698690

(9) DCB #2 (S)

10.05min 0.022ug/ml m

response 1099353

(+) = Expected Retention Time

F55864.D PBO05166.M Wed May 17 10:05:00 2006

Quantitation Report (QT Reviewed)

Signal #1 : C:\HPCHEM\2\DATA\051606-F\F55865.D\ECD1A.CH Vial: 17
Signal #2 : C:\HPCHEM\2\DATA\051606-F\F55865.D\ECD2B.CH
Acq On : 16 May 2006 8:36 pm Operator:
Sample : P4020 16/60 0.25 Inst : Instrumen
Misc : Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 17 9:54 2006 Quant Results File: PBQ05166.RES

Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 09:54:21 2006
Response via : Initial Calibration
DataAcq Meth : PSTQUICK.M

5.17.06 5.17.06

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film

Compound	RT#1	RT#2	Resp#1	Resp#2	ug/ml	ug/ml
System Monitoring Compounds						
1) S TCX	2.64	2.73	1837346	2057065	0.058	0.058
Spiked Amount	0.100	Range 46 - 122		Recovery =	58.00%	58.00%
9) S DCB	10.40	10.04	1671904	2550743	0.054	0.061
Spiked Amount	0.100	Range 26 - 86		Recovery =	54.00%	61.00%
Target Compounds						
2) L1 Ar1016 A	3.44	3.37	469862	514274	0.296	0.300
3) L1 Ar1016 B	3.93	3.87	885341	1150893	0.293	0.299
4) L1 Ar1016 C	4.09	4.03	365795	438601	0.294	0.304
Sum Ar1016 A			1720998	2103768	0.884	0.903
Average Ar1016 A					0.295	0.301
5) L2 Ar1260 A	6.67	6.35	562785	744547	0.288	0.321
6) L2 Ar1260 B	7.03	6.82	670680	1091231	0.285	0.319
7) L2 Ar1260 C	7.92	7.53	510810	774876	0.279	0.318
8) L2 Ar1260 D	8.49	8.59	500733	1688943	0.278	0.299
Sum Ar1260 A			2245008	4299597	1.130	1.256
Average Ar1260 A					0.283	0.314

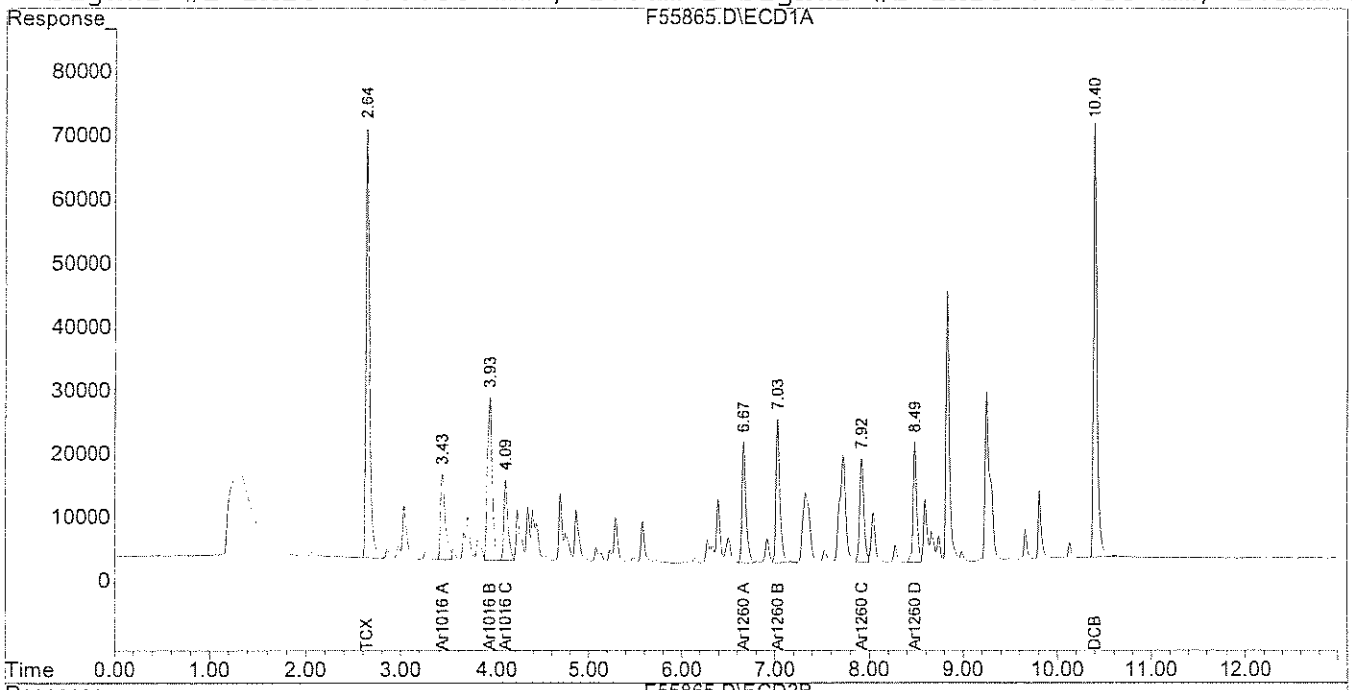
(f)=RT Delta > 1/2 Window (#)=Amounts differ by > 25% (m)=manual int.

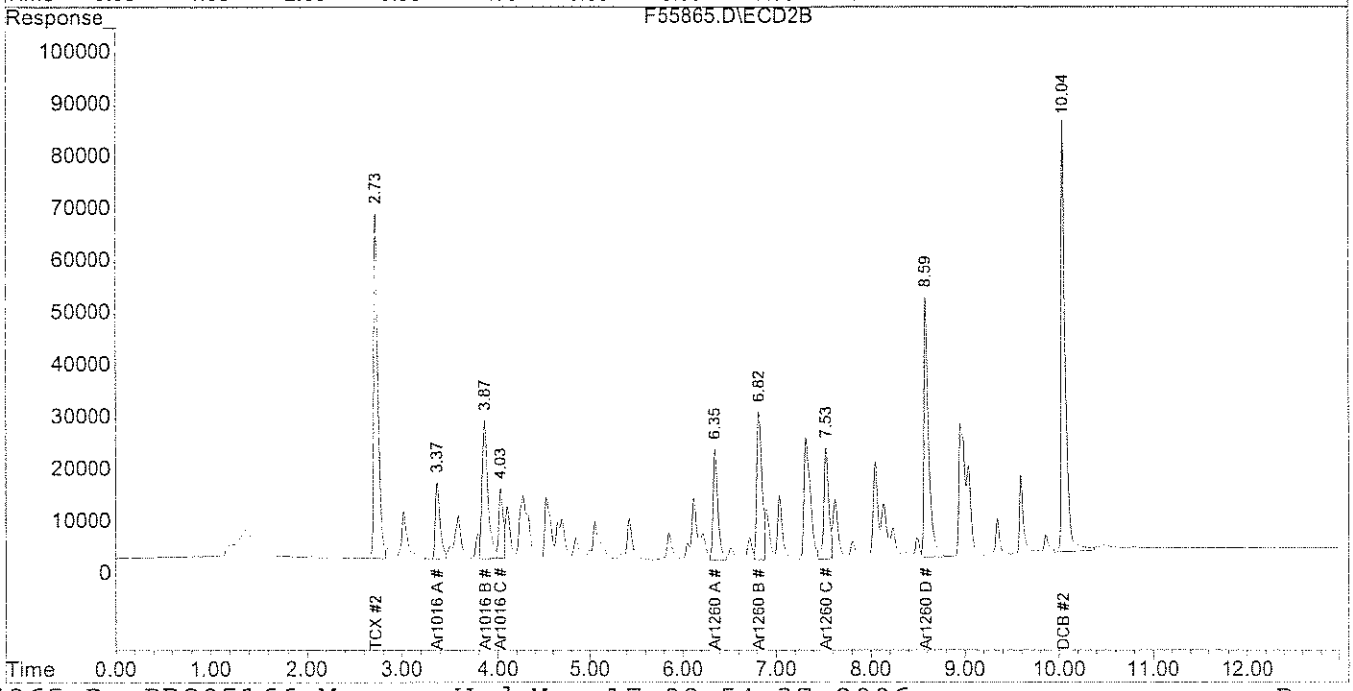
Quantitation Report (QT Reviewed)

Signal #1 : C:\HPCHEM\2\DATA\051606-F\F55865.D\ECD1A.CH Vial: 17
Signal #2 : C:\HPCHEM\2\DATA\051606-F\F55865.D\ECD2B.CH
Acq On : 16 May 2006 8:36 pm Operator:
Sample : P4020 16/60 0.25 Inst : Instrumen
Misc : Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 17 9:54 2006 Quant Results File: PBQ05166.RES

Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 09:54:21 2006
Response via : Multiple Level Calibration
DataAcq Meth : PSTQUICK.M

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film





Quantitation Report (QT Reviewed)

Signal #1 : C:\HPCHEM\2\DATA\051606-F\F55866.D\ECD1A.CH Vial: 18
Signal #2 : C:\HPCHEM\2\DATA\051606-F\F55866.D\ECD2B.CH
Acq On : 16 May 2006 8:51 pm Operator:
Sample : P4021 16/60 0.5 Inst : Instrumen
Misc : Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 17 10:05 2006 Quant Results File: PBQ05166.RES

Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 09:54:21 2006
Response via : Initial Calibration
DataAcq Meth : PSTQUICK.M

5.17.06 5.17.06

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film

Compound	RT#1	RT#2	Resp#1	Resp#2	ug/ml	ug/ml
System Monitoring Compounds						
1) S TCX	2.64	2.73	3525739	3853418	0.112	0.110
Spiked Amount 0.100	Range 46 - 122		Recovery =		112.00%	110.00%
9) S DCB	10.39	10.04	3091651	4358733	0.104	0.108
Spiked Amount 0.100	Range 26 - 86		Recovery =		104.00%#	108.00%#
Target Compounds						
2) L1 Ar1016 A	3.43	3.37	867015	920329	0.561	0.558
3) L1 Ar1016 B	3.93	3.87	1650244	2072594	0.559	0.557
4) L1 Ar1016 C	4.09	4.03	686360	787671	0.560	0.564
Sum Ar1016 A			3203619	3780595	1.680	1.679
Average Ar1016 A					0.560	0.560
5) L2 Ar1260 A	6.67	6.35	1071560	1271657	0.562	0.575
6) L2 Ar1260 B	7.03	6.82	1268608	1900394	0.552	0.579
7) L2 Ar1260 C	7.92	7.53	993398	1384156	0.549	0.587
8) L2 Ar1260 D	8.49	8.59	972866	3155830	0.547	0.571
Sum Ar1260 A			4306432	7712037	2.210	2.312
Average Ar1260 A					0.553	0.578

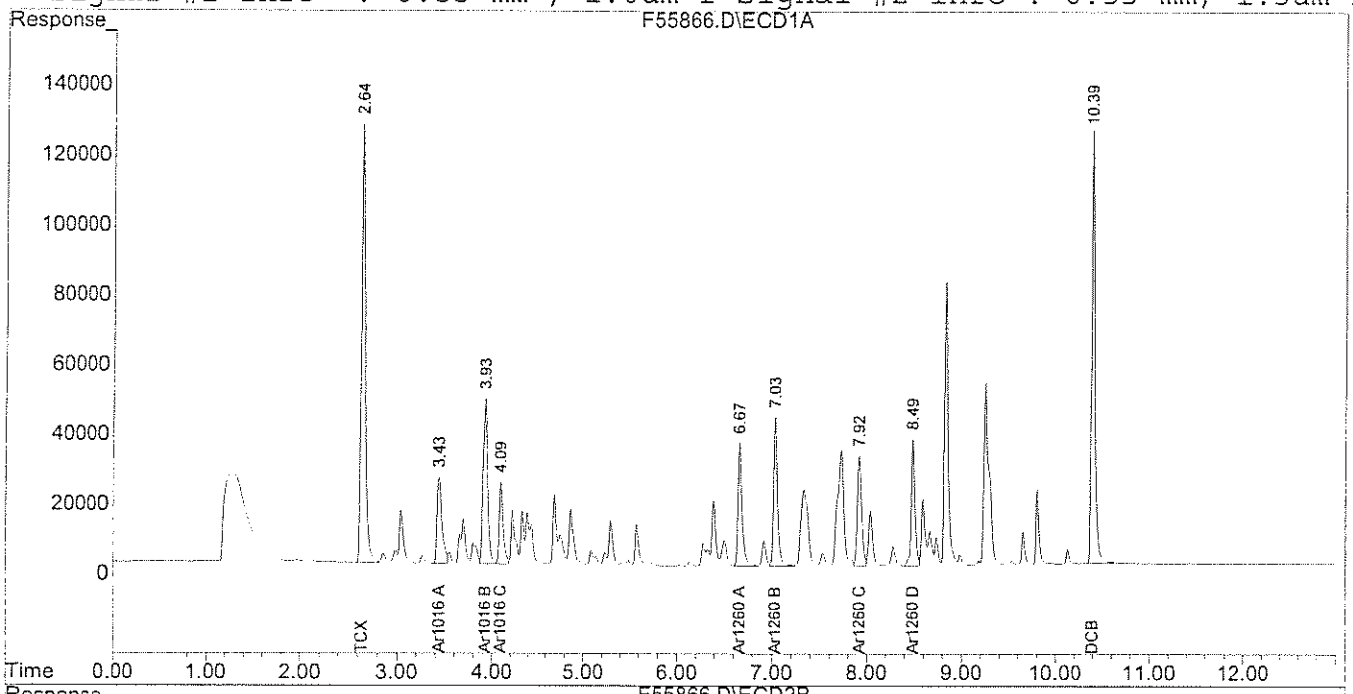
(f)=RT Delta > 1/2 Window (#)=Amounts differ by > 25% (m)=manual int

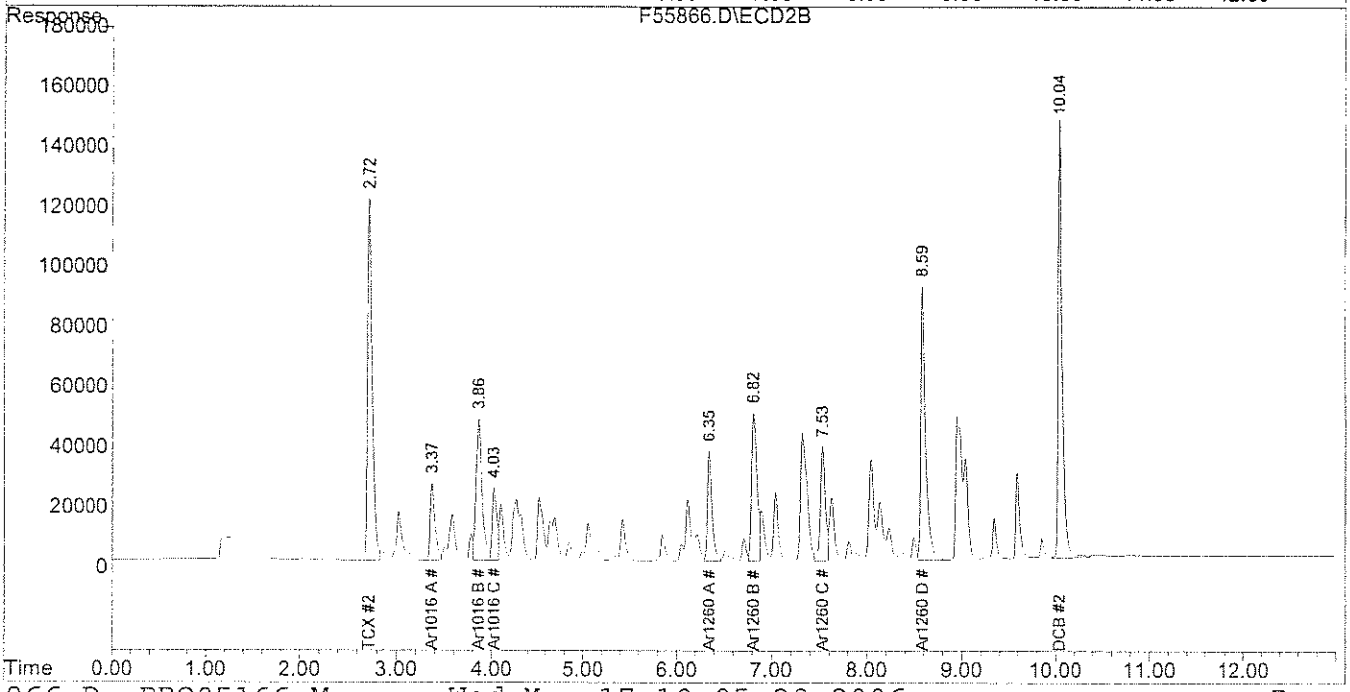
Quantitation Report (QT Reviewed)

Signal #1 : C:\HPCHEM\2\DATA\051606-F\F55866.D\ECD1A.CH Vial: 18
Signal #2 : C:\HPCHEM\2\DATA\051606-F\F55866.D\ECD2B.CH
Acq On : 16 May 2006 8:51 pm Operator:
Sample : P4021 16/60 0.5 Inst : Instrumen
Misc : Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 17 10:05 2006 Quant Results File: PBQ05166.RES

Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 09:54:21 2006
Response via : Multiple Level Calibration
DataAcq Meth : PSTQUICK.M

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film





Quantitation Report (QT Reviewed)

Signal #1 : C:\HPCHEM\2\DATA\051606-F\F55867.D\ECD1A.CH Vial: 19
Signal #2 : C:\HPCHEM\2\DATA\051606-F\F55867.D\ECD2B.CH
Acq On : 16 May 2006 9:20 pm Operator:
Sample : P4022 16/60 1.0 Inst : Instrumen
Misc : Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 17 10:05 2006 Quant Results File: PBQ05166.RES

Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 09:54:21 2006
Response via : Initial Calibration
DataAcq Meth : PSTQUICK.M

5.17.06 5.17.06

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film

Compound	RT#1	RT#2	Resp#1	Resp#2	ug/ml	ug/ml
System Monitoring Compounds						
1) S TCX	2.64	2.73	6920671	7382493	0.219	0.211
Spiked Amount	0.100	Range 46 - 122		Recovery =	219.00%#	211.00%#
9) S DCB	10.40	10.04	5840708	7809116	0.200	0.200
Spiked Amount	0.100	Range 26 - 86		Recovery =	200.00%#	200.00%#
Target Compounds						
2) L1 Ar1016 A	3.43	3.37	1663419	1708138	1.092	1.059
3) L1 Ar1016 B	3.93	3.87	3193003	3862796	1.095	1.058
4) L1 Ar1016 C	4.09	4.03	1338659	1462474	1.101	1.068
Sum Ar1016 A			6195082	7033407	3.288	3.185
Average Ar1016 A					1.096	1.062
5) L2 Ar1260 A	6.67	6.35	2047135	2313020	1.087	1.076
6) L2 Ar1260 B	7.03	6.82	2456273	3457427	1.083	1.080
7) L2 Ar1260 C	7.92	7.53	1964211	2559137	1.092	1.107
8) L2 Ar1260 D	8.49	8.59	1930282	5977809	1.093	1.094
Sum Ar1260 A			8397902	14307393	4.355	4.357
Average Ar1260 A					1.089	1.089

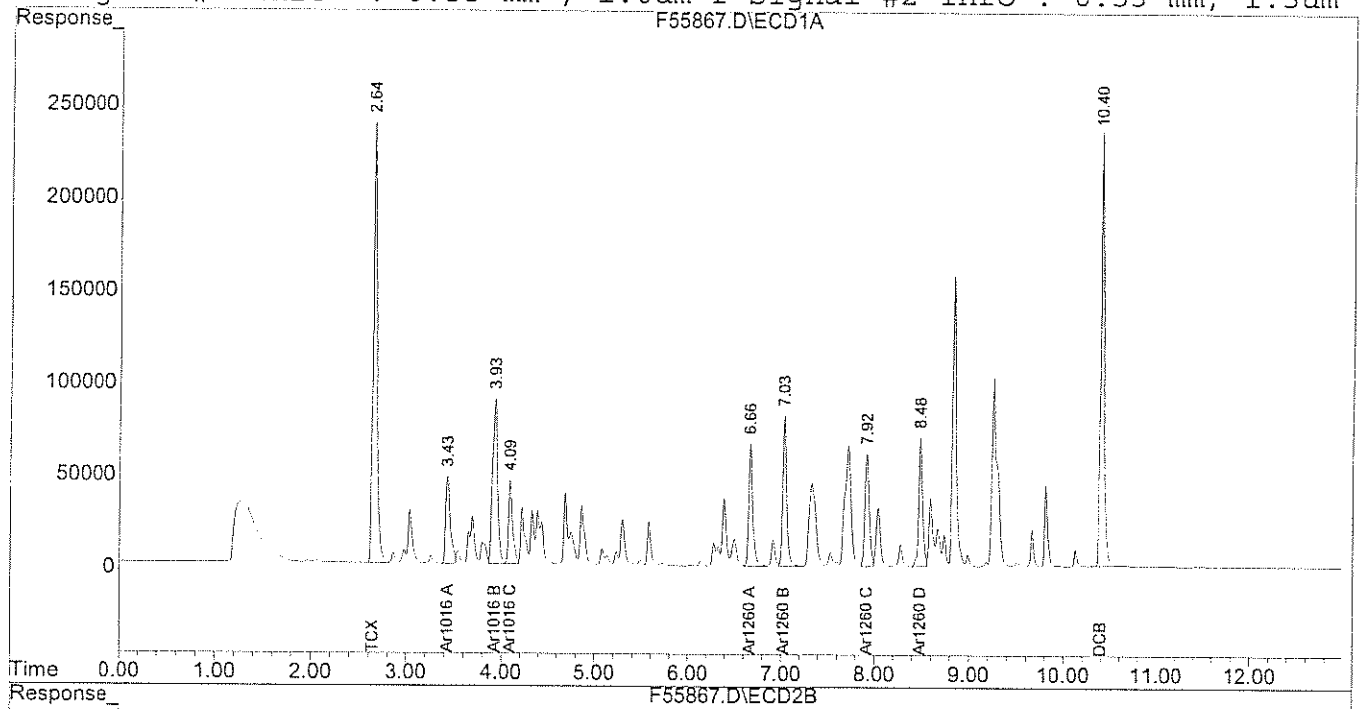
(f)=RT Delta > 1/2 Window (#)=Amounts differ by > 25% (m)=manual int.

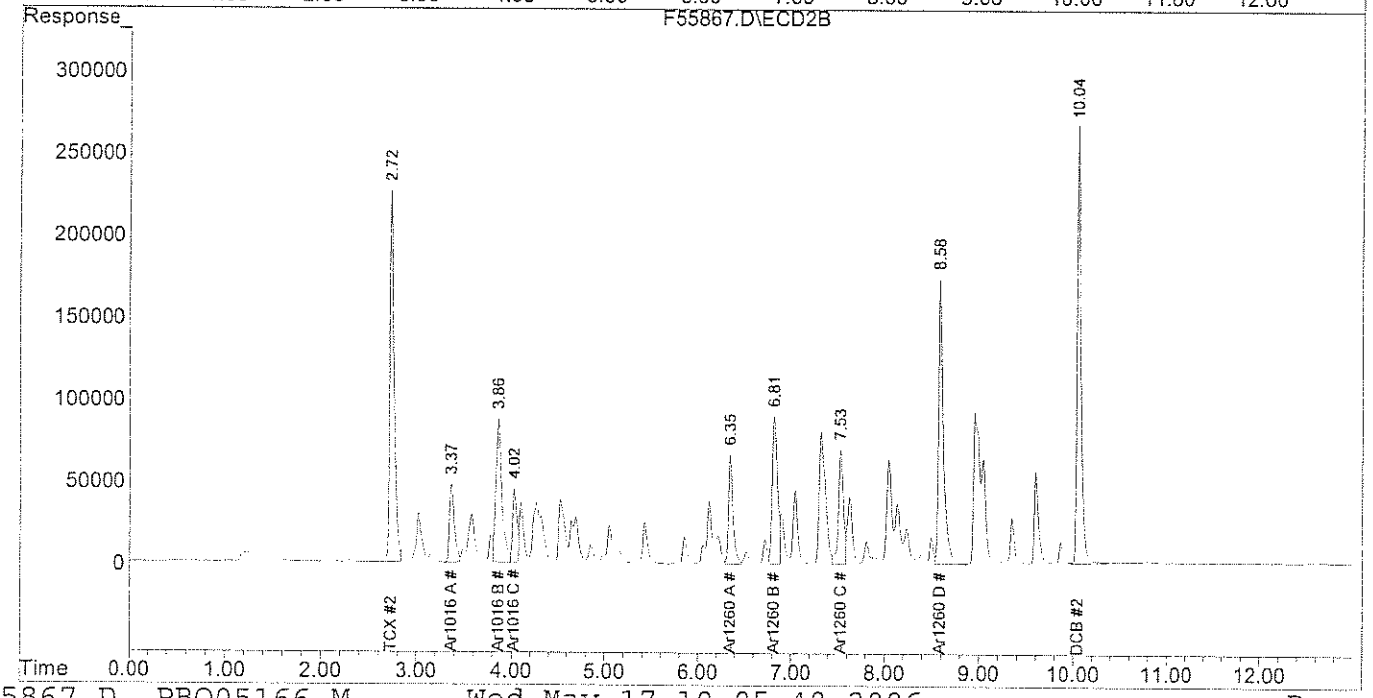
Quantitation Report (QT Reviewed)

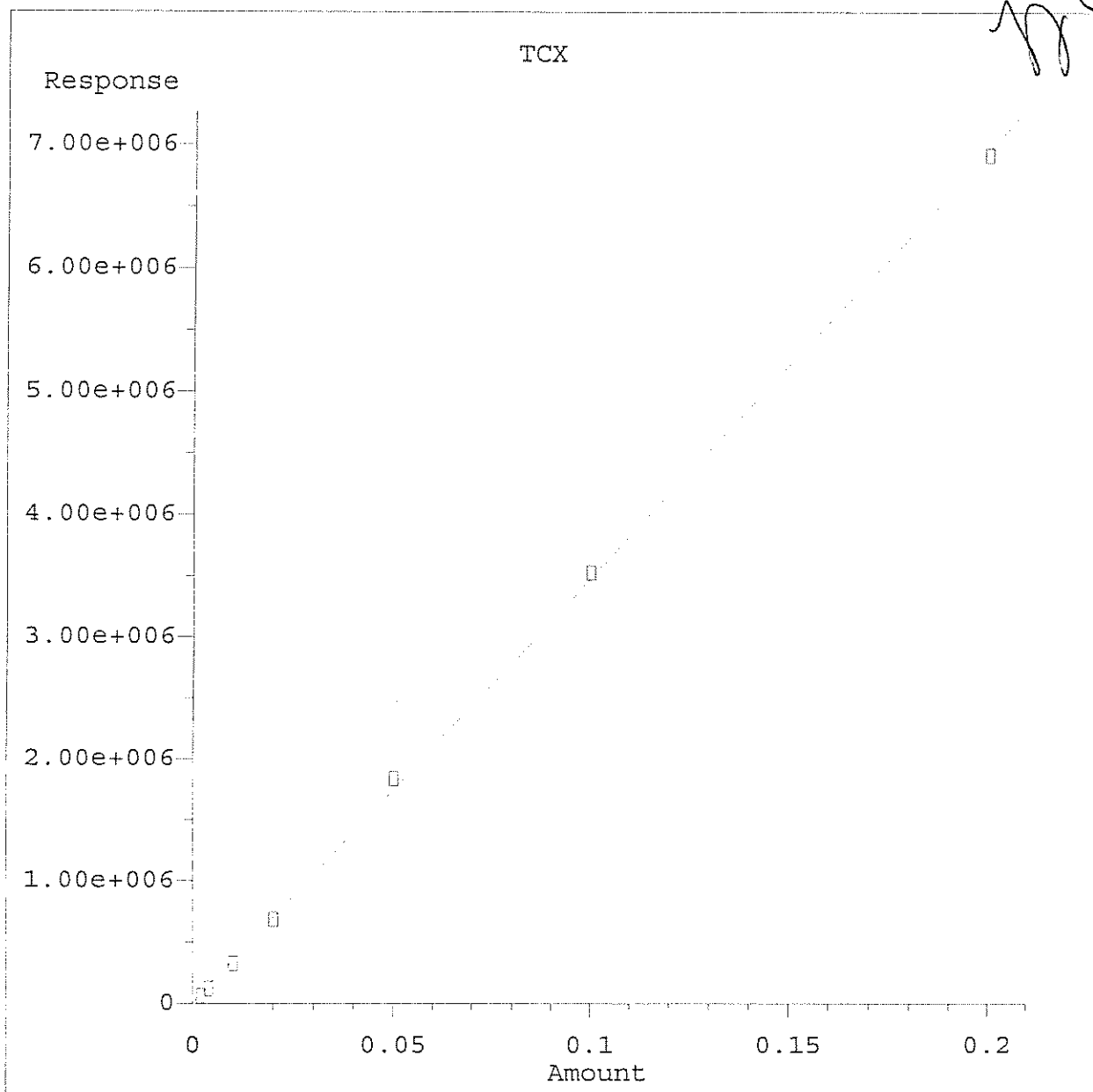
Signal #1 : C:\HPCHEM\2\DATA\051606-F\F55867.D\ECD1A.CH Vial: 19
Signal #2 : C:\HPCHEM\2\DATA\051606-F\F55867.D\ECD2B.CH
Acq On : 16 May 2006 9:20 pm Operator:
Sample : P4022 16/60 1.0 Inst : Instrumen
Misc : Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 17 10:05 2006 Quant Results File: PBQ05166.RES

Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 09:54:21 2006
Response via : Multiple Level Calibration
DataAcq Meth : PSTQUICK.M

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film



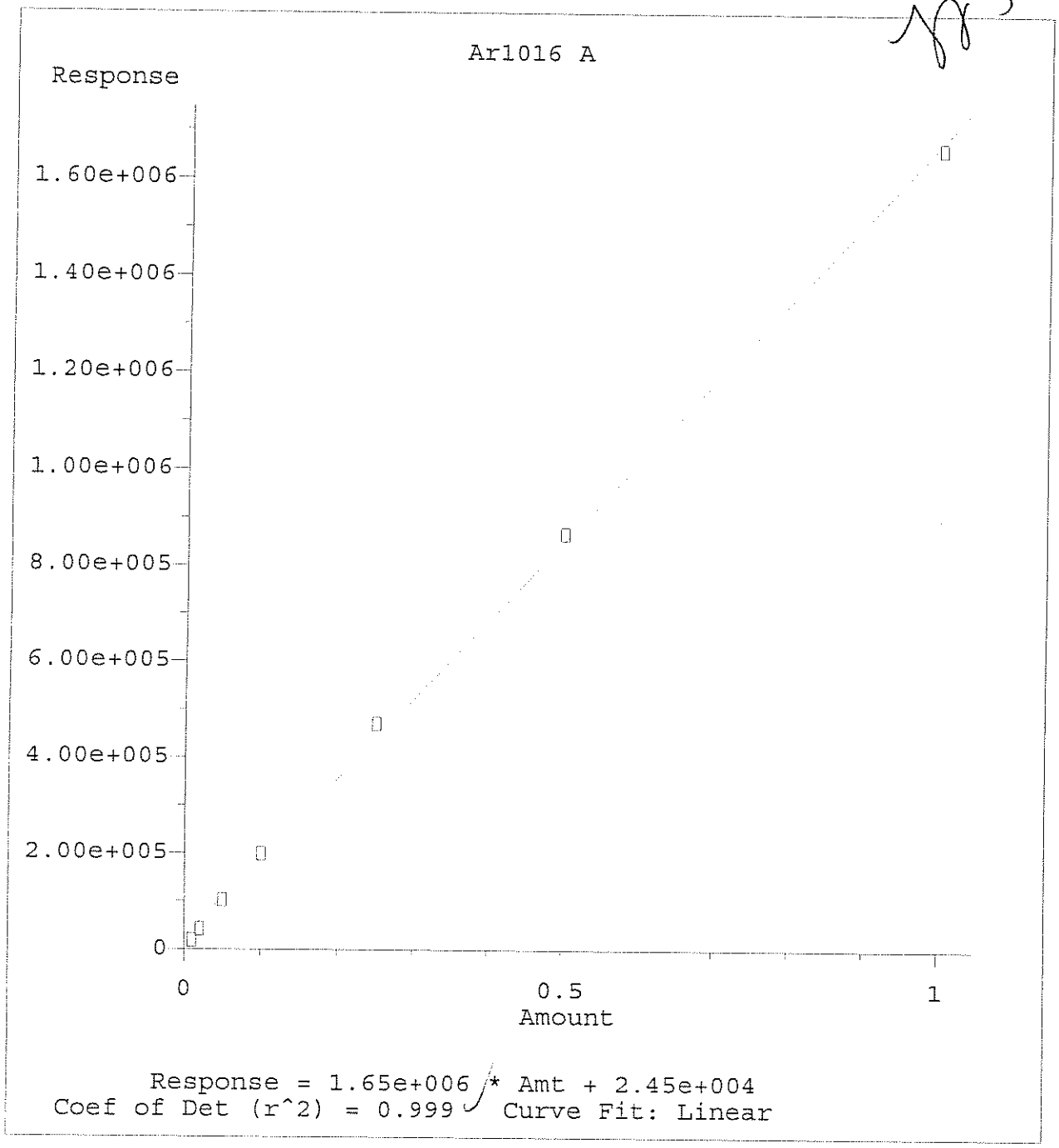




Response = 3.47e+007 * Amt + 1.04e+004
Coef of Det (r^2) = 1.000 Curve Fit: Linear

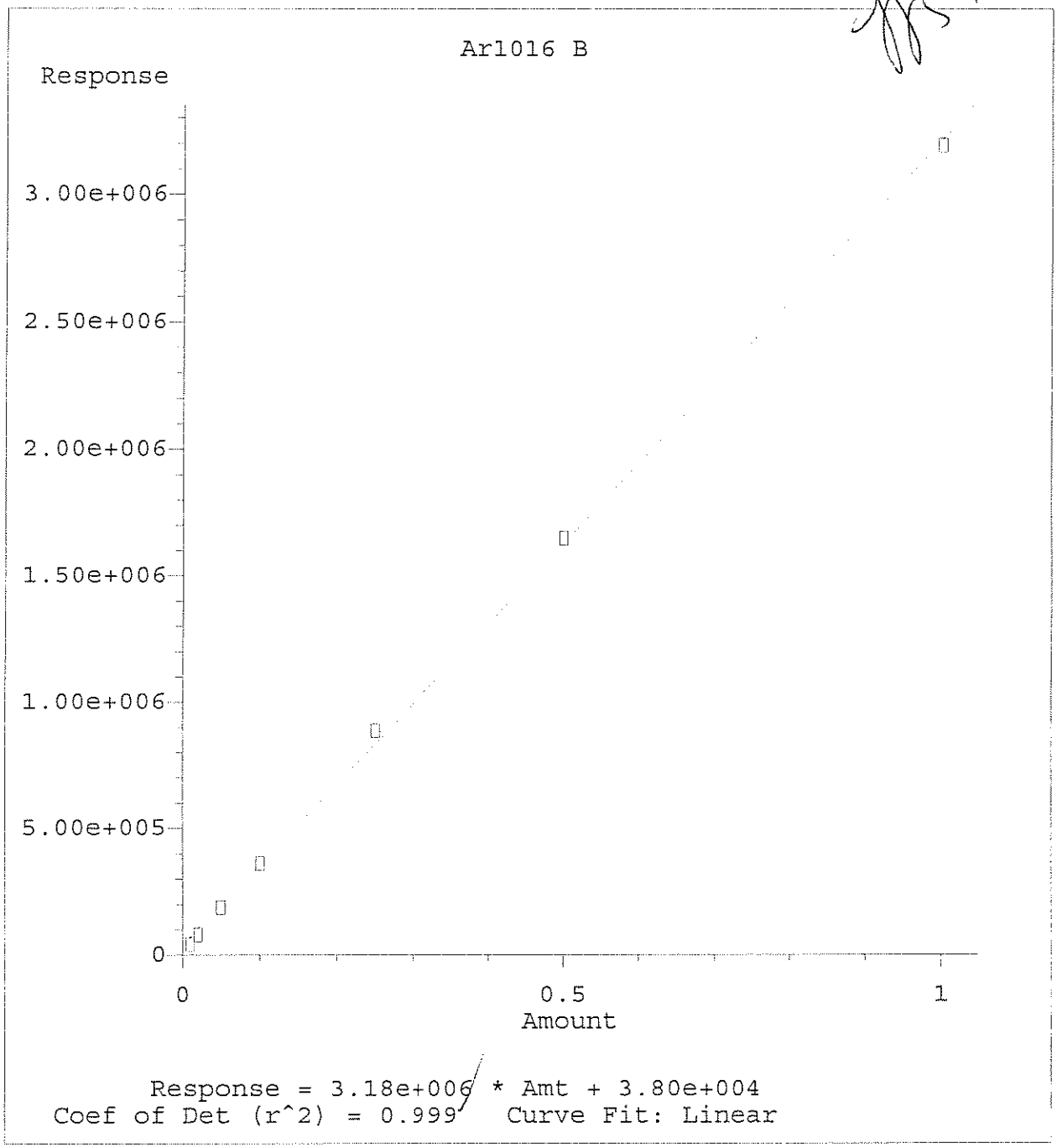
Method Name: C:\HPCHEM\2\METHODS\PBQ05166.M
Calibration Table Last Updated: Wed May 17 10:06:32 2006

5-17-06



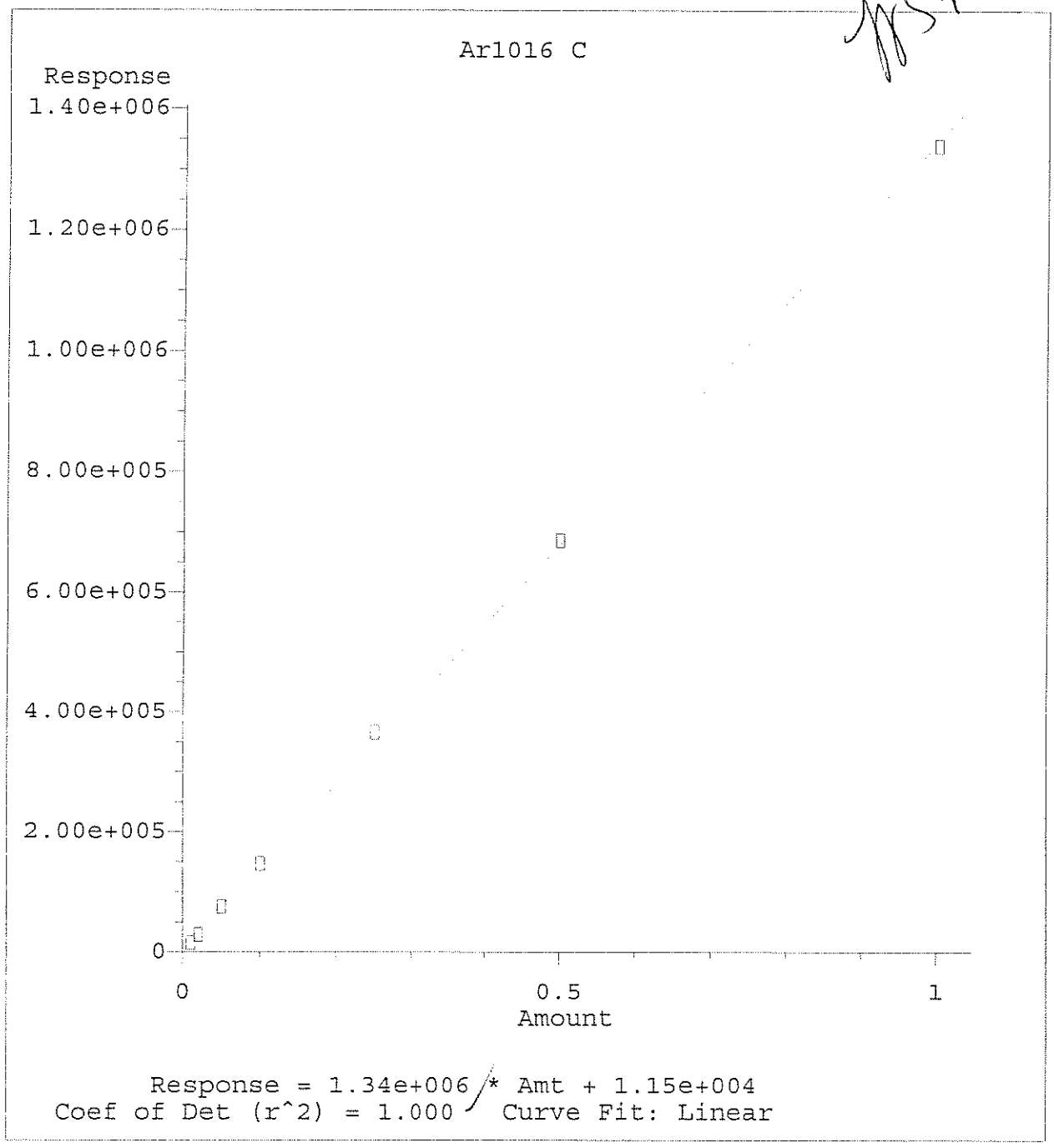
Method Name: C:\HPCHEM\2\METHODS\PBQ05166.M
Calibration Table Last Updated: Wed May 17 10:06:32 2006

5-17-06



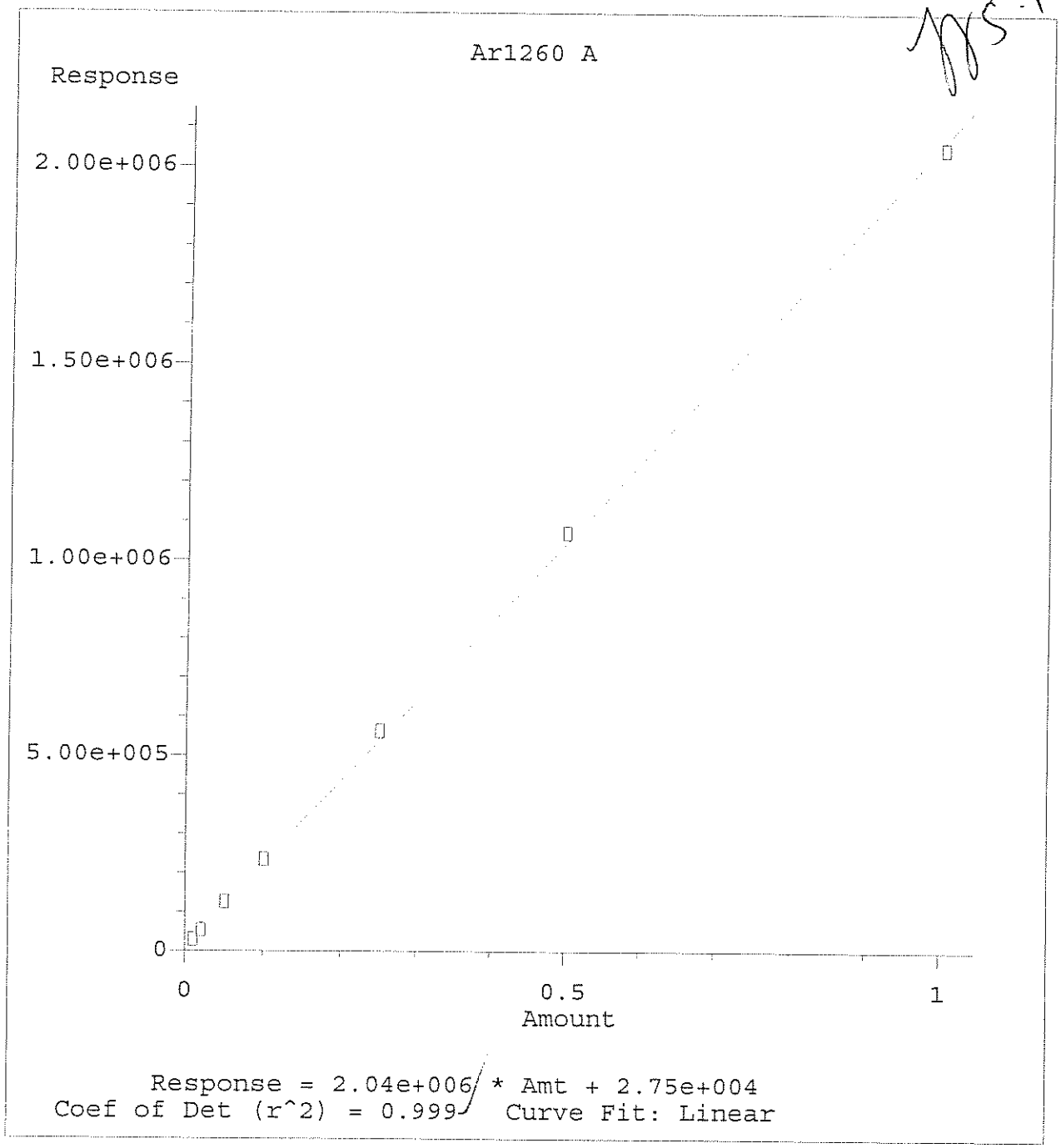
Method Name: C:\HPCHEM\2\METHODS\PBQ05166.M
Calibration Table Last Updated: Wed May 17 10:06:32 2006

5.17.06



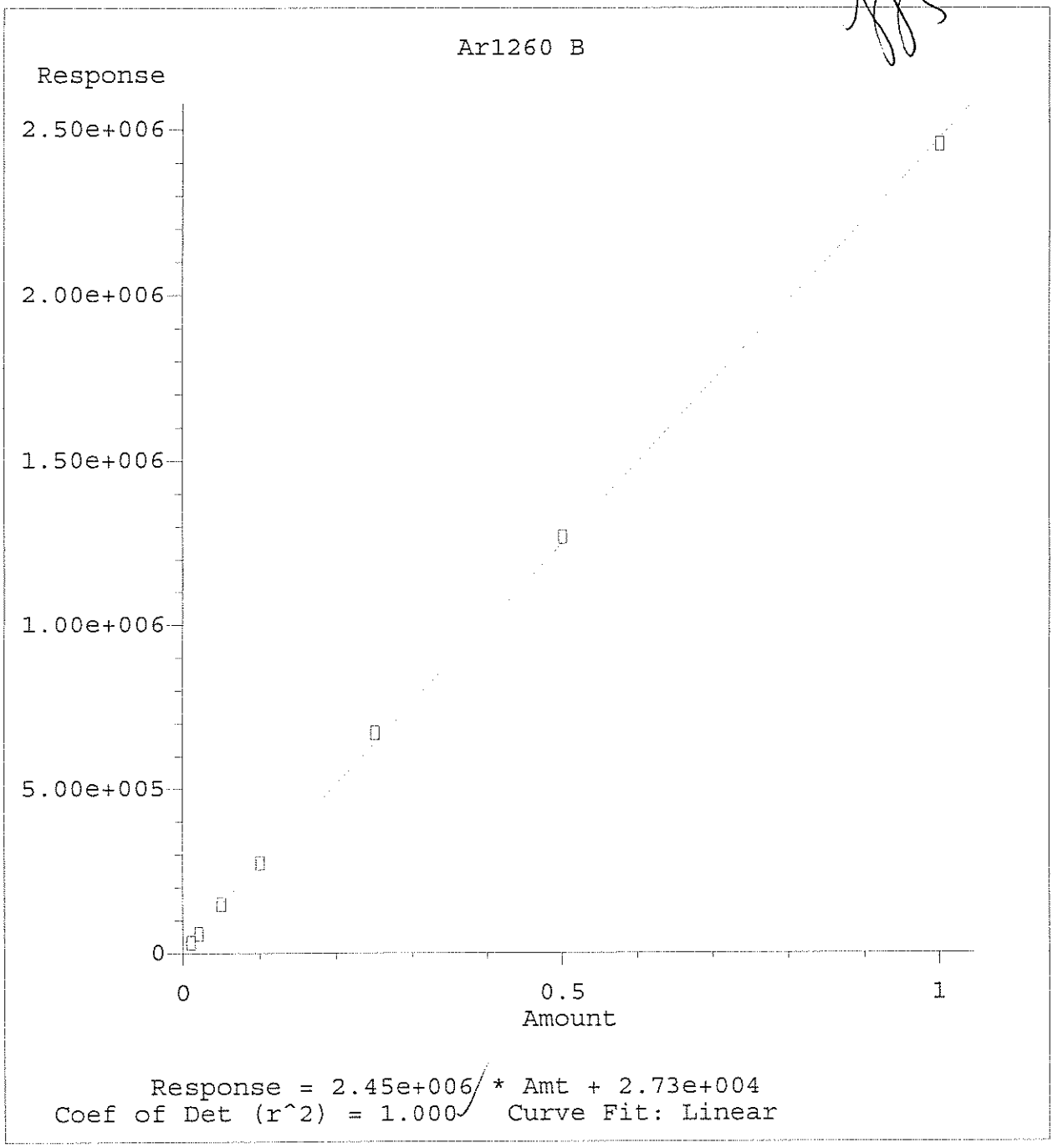
Method Name: C:\HPCHEM\2\METHODS\PBQ05166.M
Calibration Table Last Updated: Wed May 17 10:06:32 2006

5.17.06



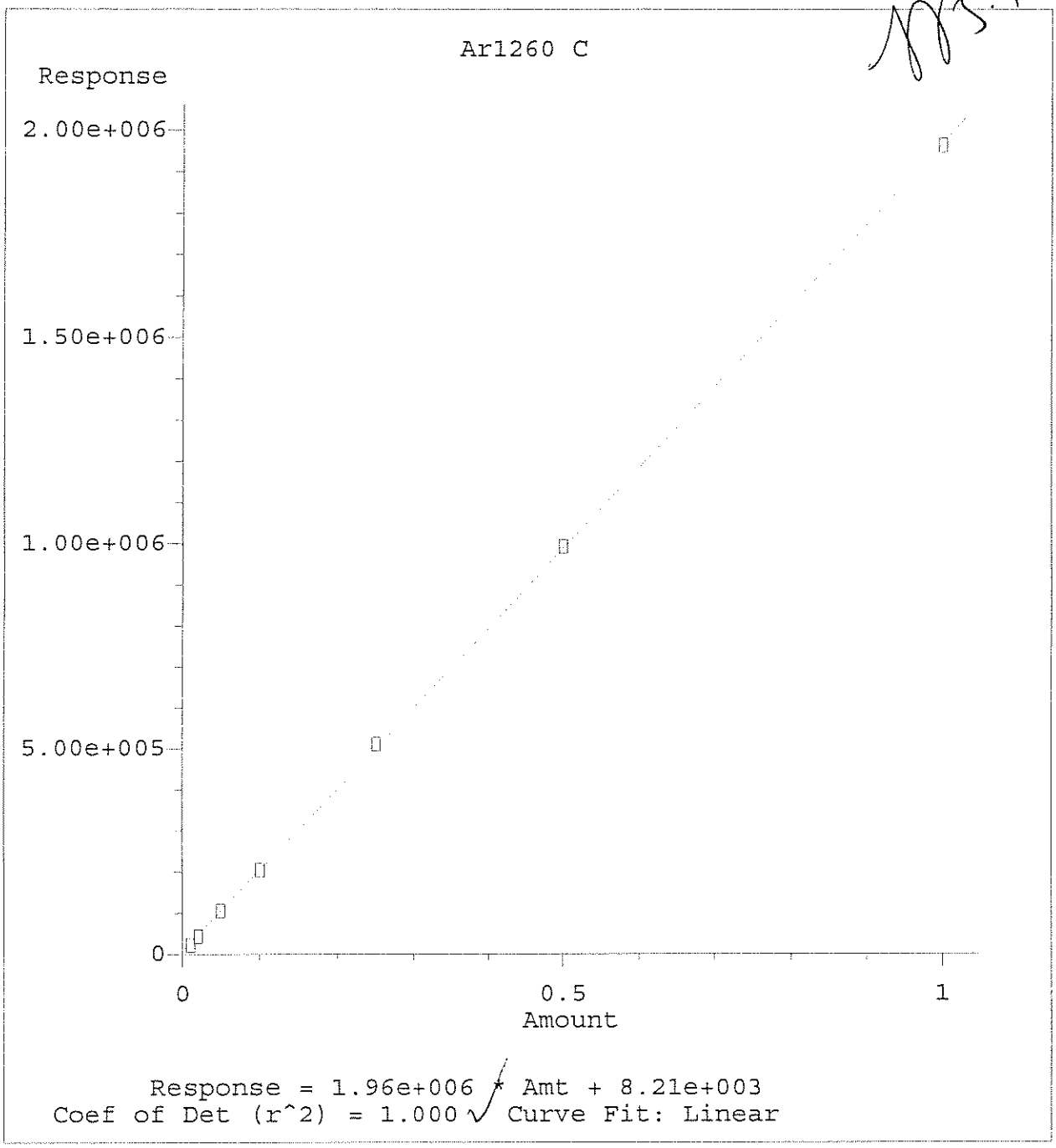
Method Name: C:\HPCHEM\2\METHODS\PBQ05166.M
Calibration Table Last Updated: Wed May 17 10:06:32 2006

5.17.06

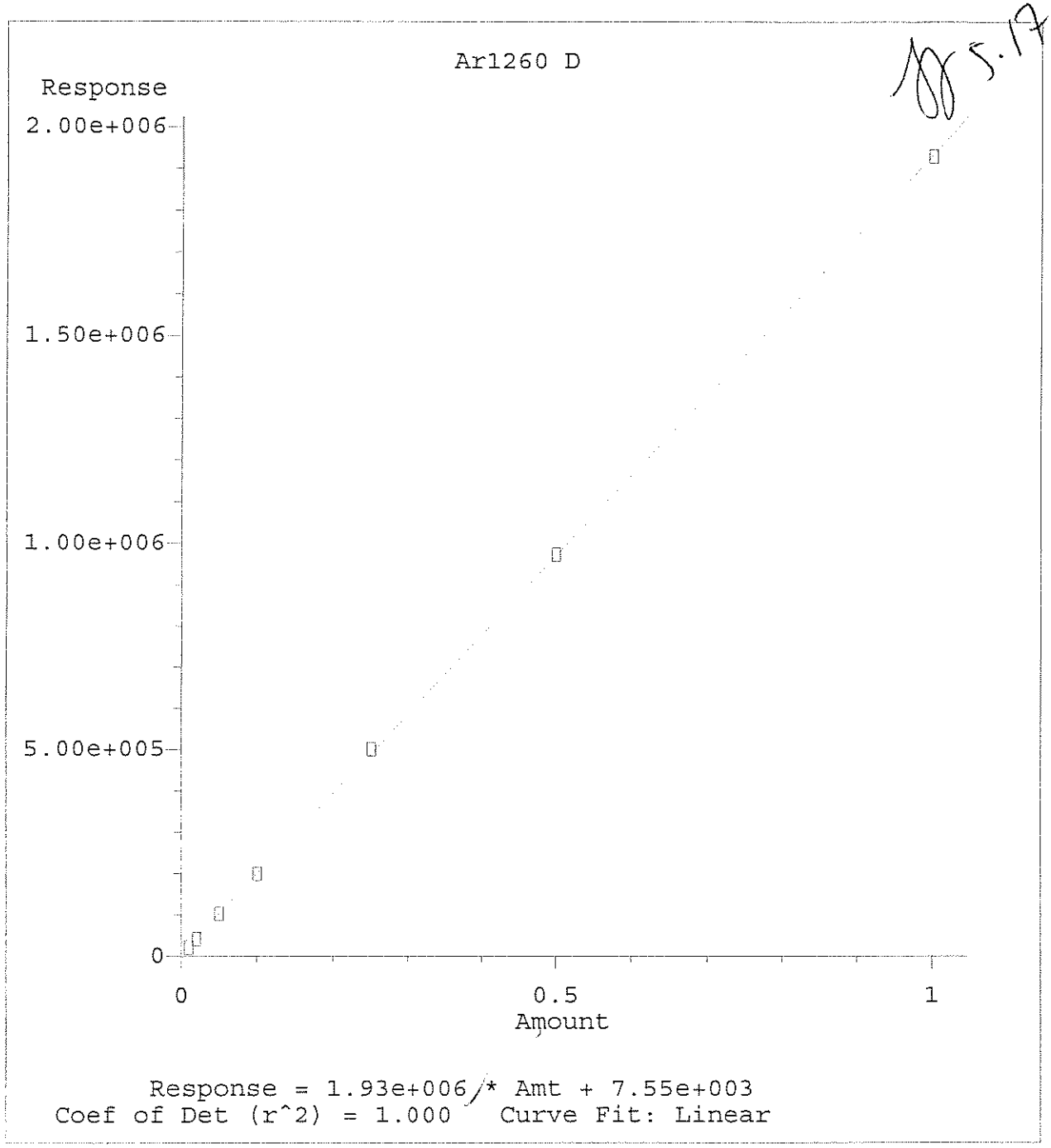


Method Name: C:\HPCHEM\2\METHODS\PBQ05166.M
Calibration Table Last Updated: Wed May 17 10:06:32 2006

5.17.06



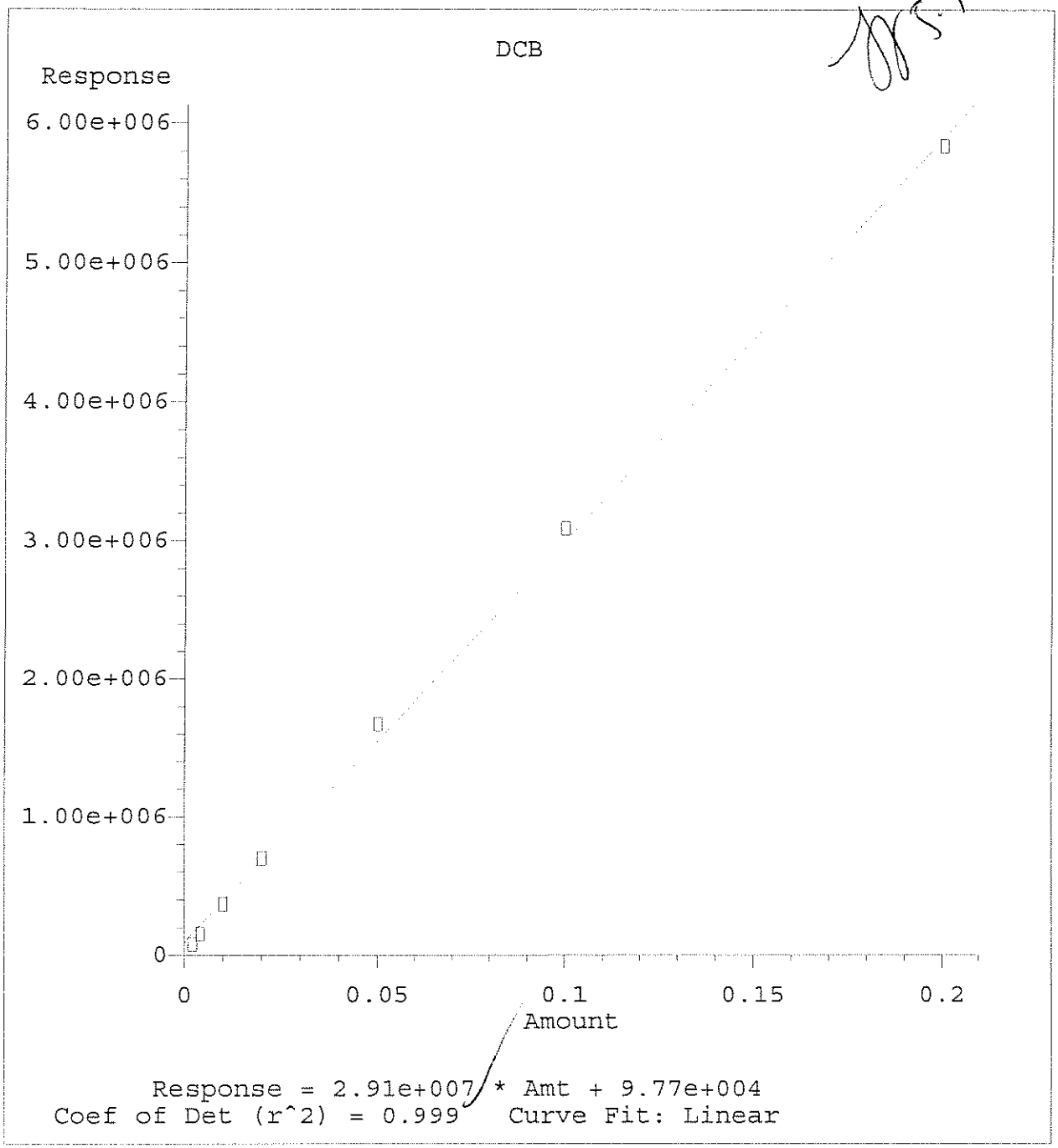
Method Name: C:\HPCHEM\2\METHODS\PBQ05166.M
Calibration Table Last Updated: Wed May 17 10:06:32 2006



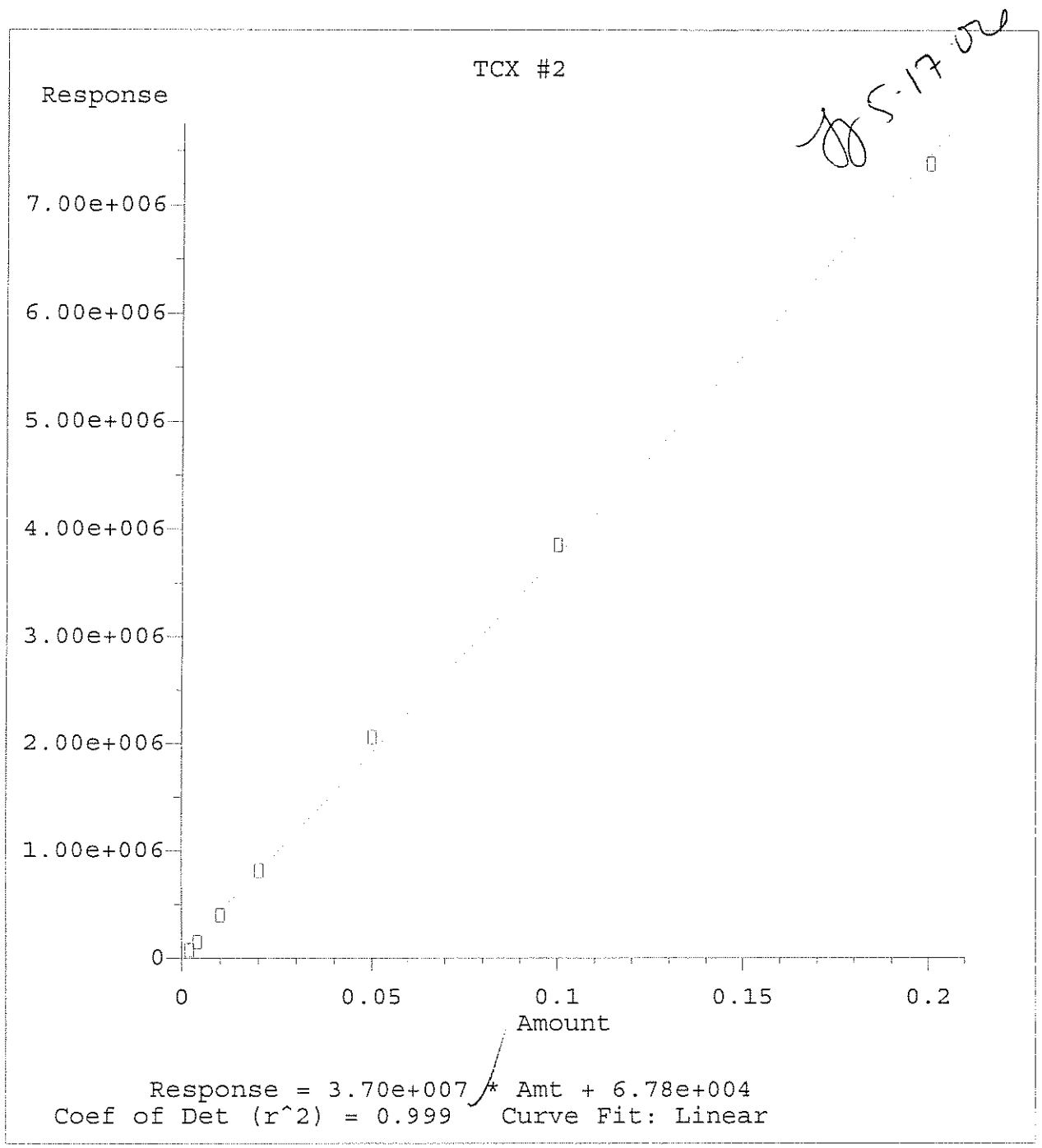
5.17.06

Method Name: C:\HPCHEM\2\METHODS\PBQ05166.M
Calibration Table Last Updated: Wed May 17 10:06:32 2006

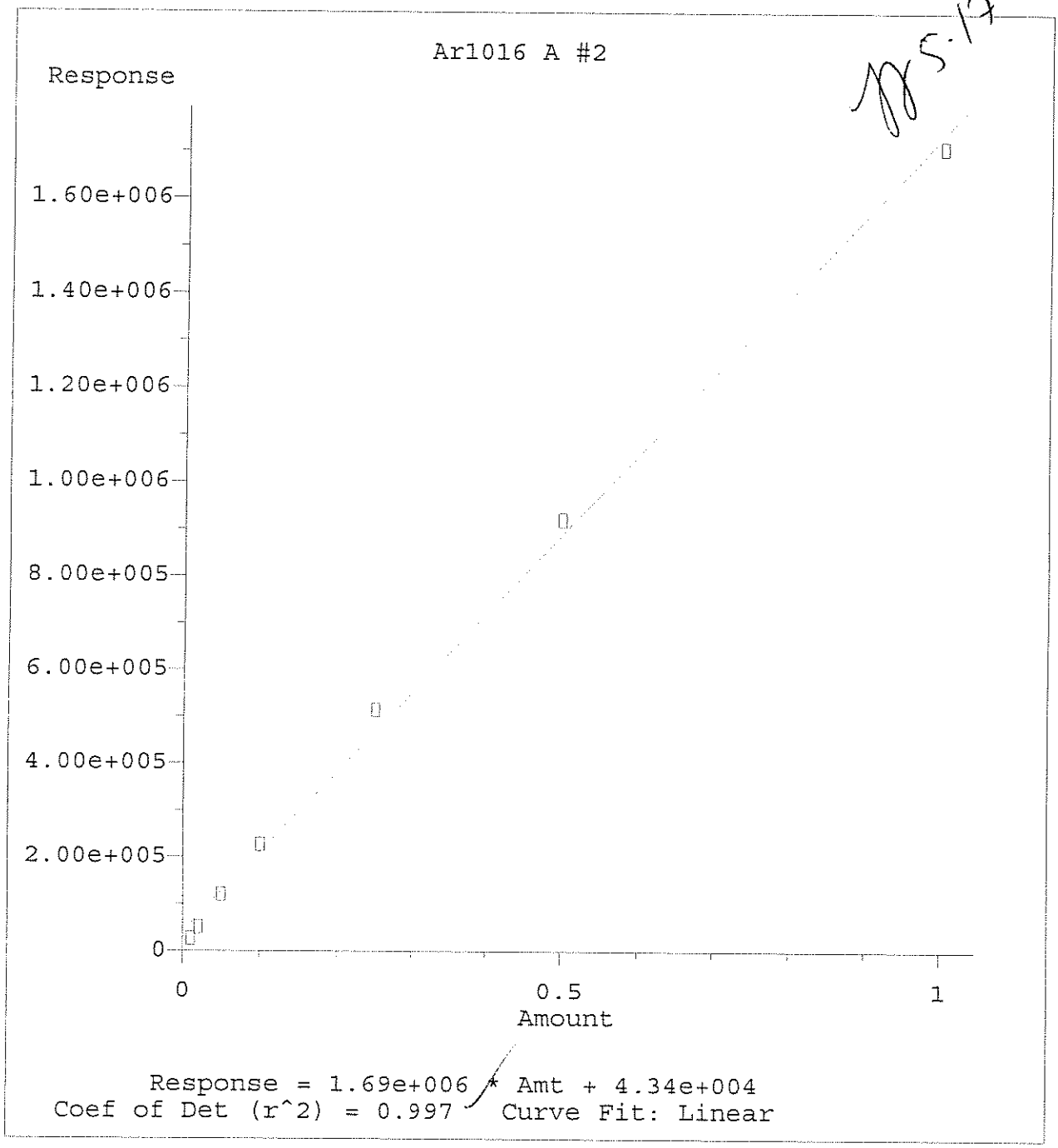
5-17-06



Method Name: C:\HPCHEM\2\METHODS\PBQ05166.M
Calibration Table Last Updated: Wed May 17 10:06:32 2006

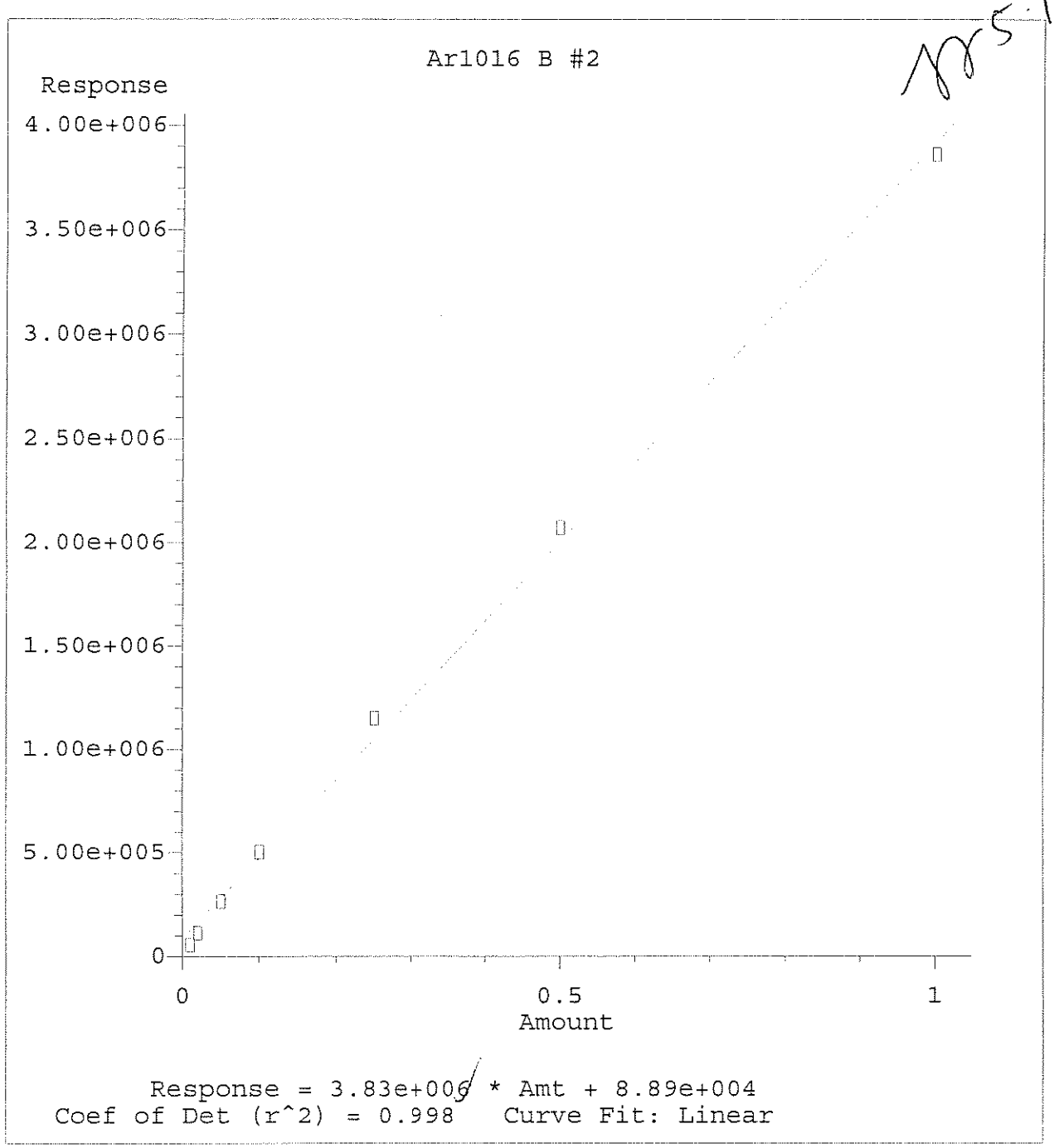


Method Name: C:\HPCHEM\2\METHODS\PBQ05166.M
Calibration Table Last Updated: Wed May 17 10:06:32 2006



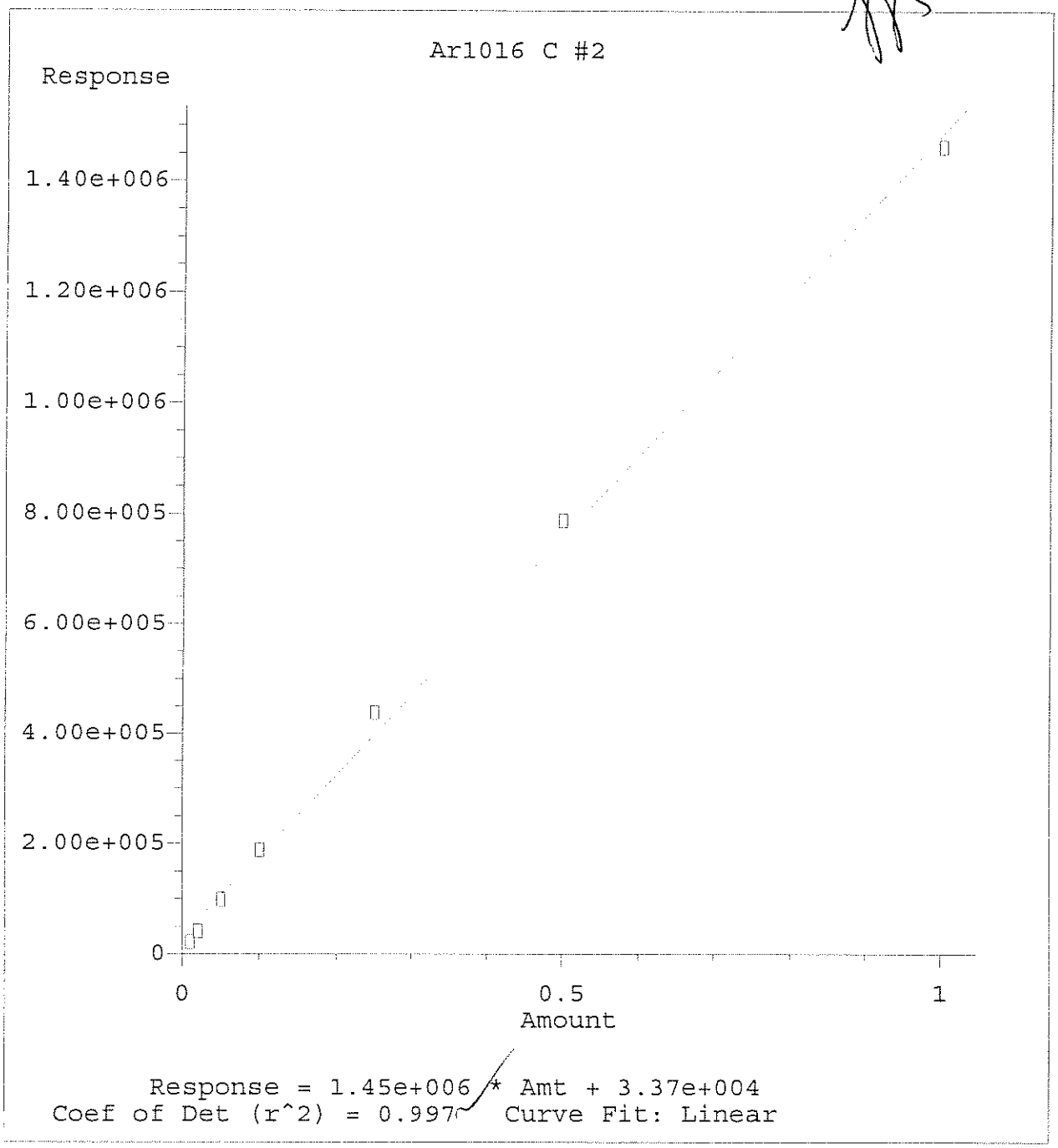
5.17.06

Method Name: C:\HPCHEM\2\METHODS\PBQ05166.M
Calibration Table Last Updated: Wed May 17 10:06:32 2006



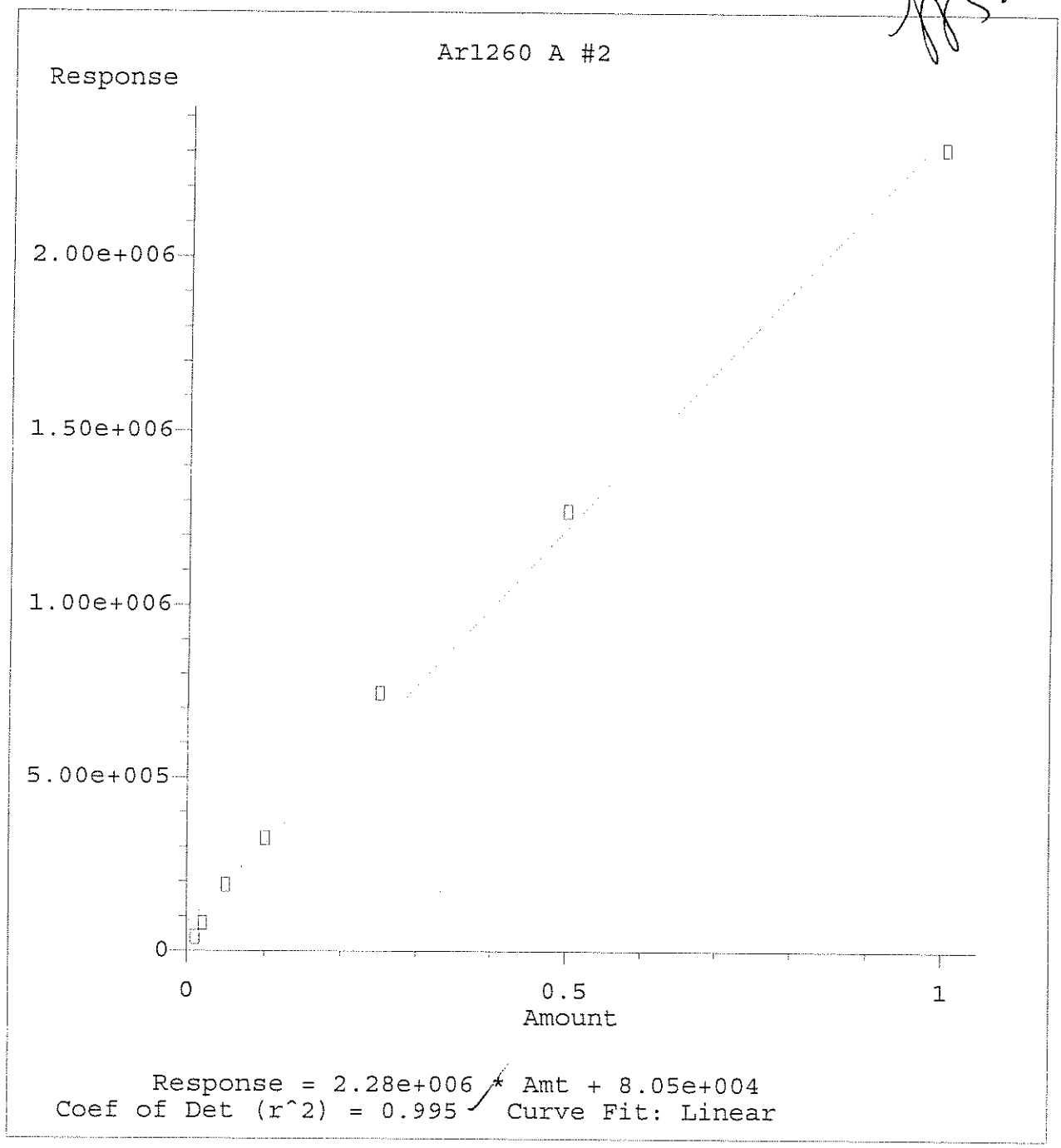
Method Name: C:\HPCHEM\2\METHODS\PBQ05166.M
Calibration Table Last Updated: Wed May 17 10:06:32 2006

5.17.06



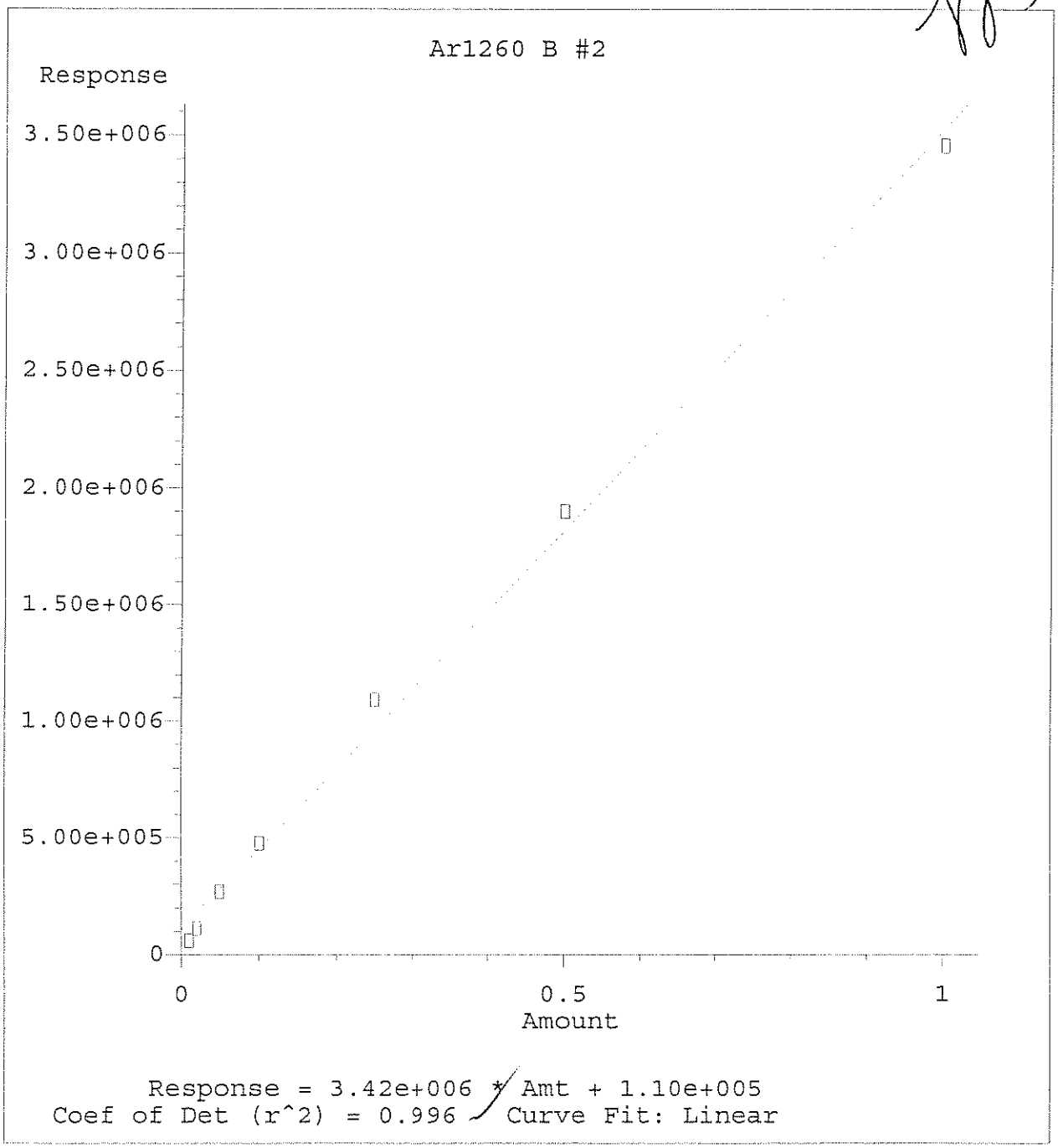
Method Name: C:\HPCHEM\2\METHODS\PBQ05166.M
Calibration Table Last Updated: Wed May 17 10:06:32 2006

5.17.06



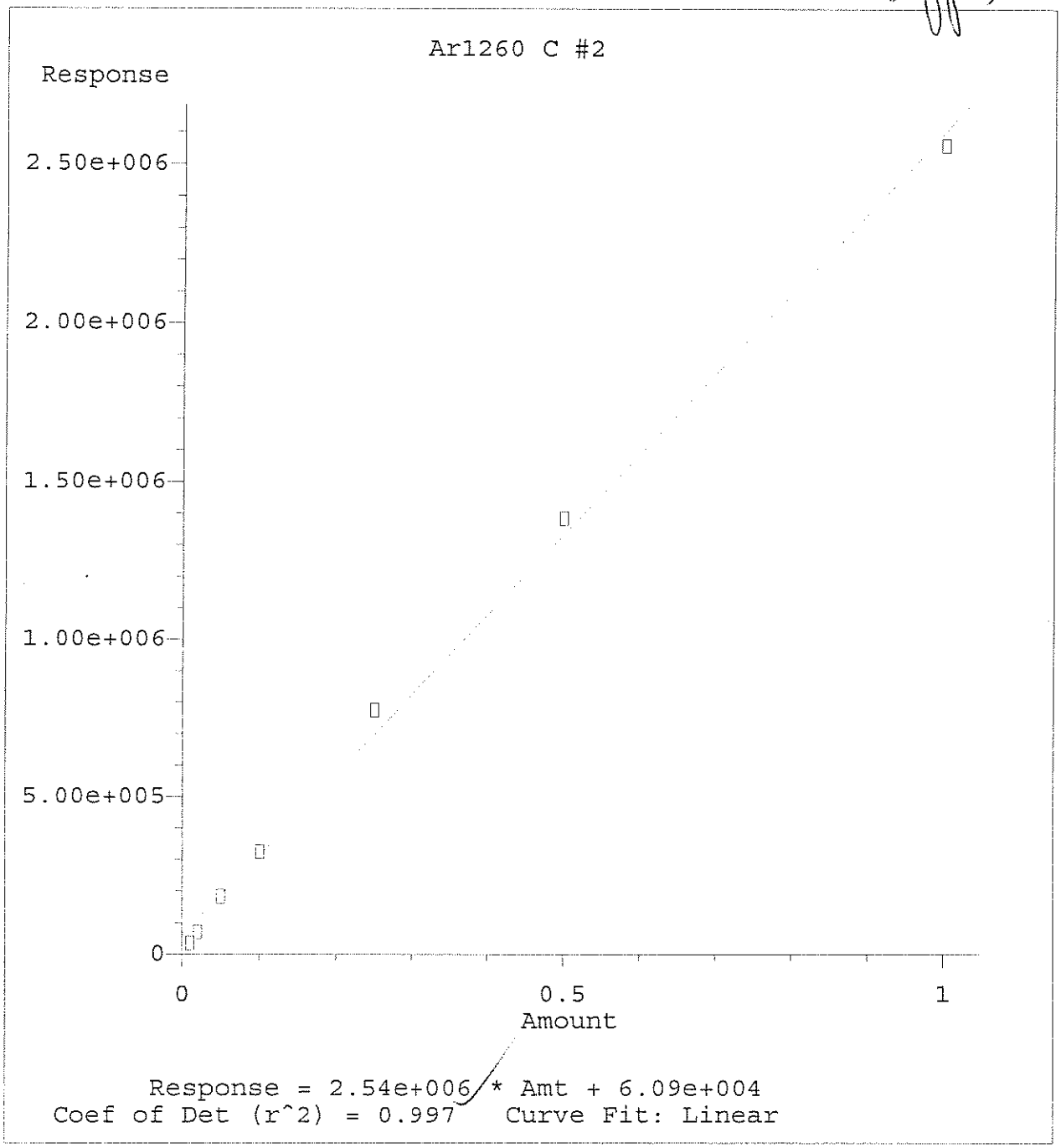
Method Name: C:\HPCHEM\2\METHODS\PBQ05166.M
Calibration Table Last Updated: Wed May 17 10:06:32 2006

5.17.06



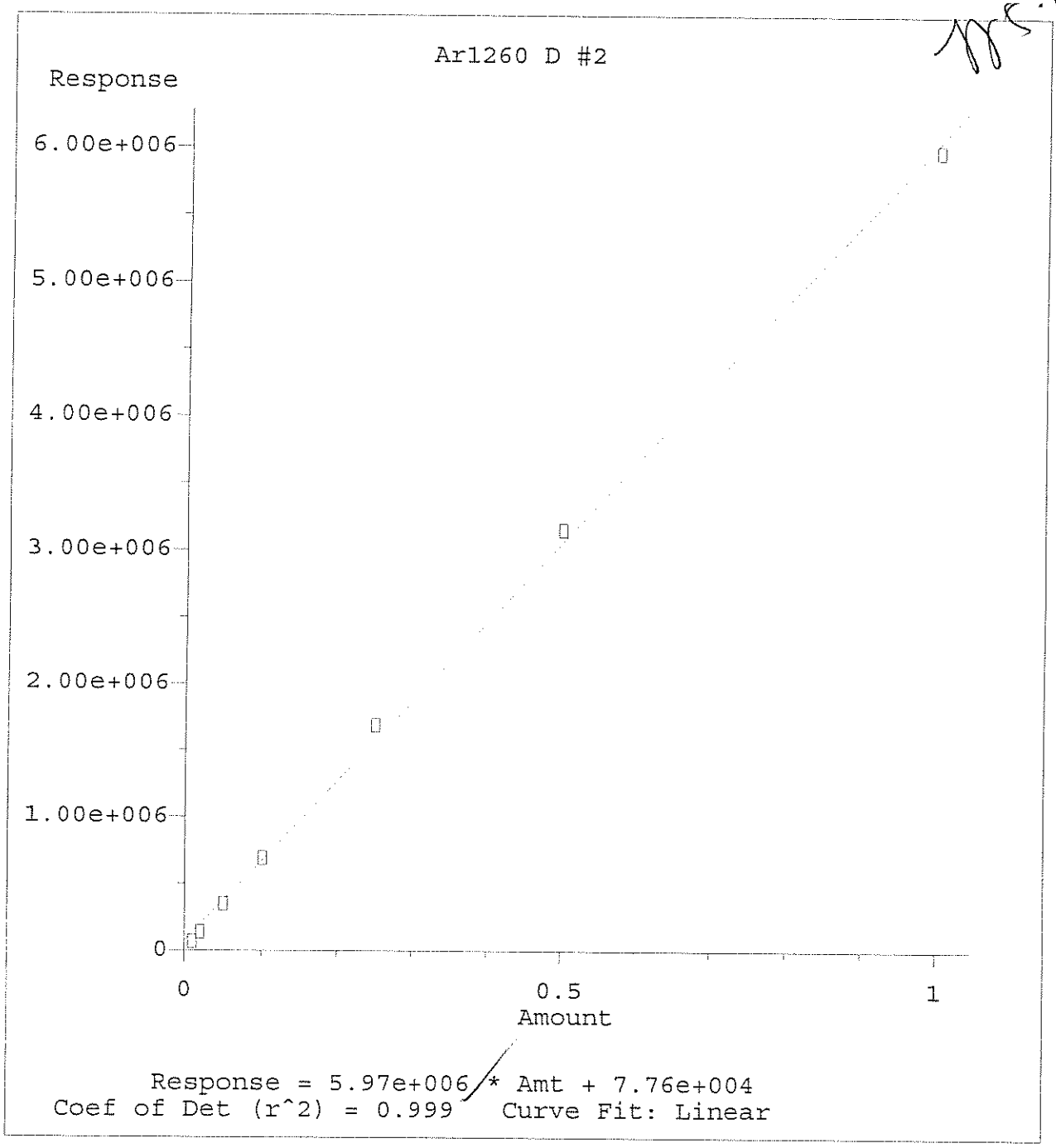
Method Name: C:\HPCHEM\2\METHODS\PBQ05166.M
Calibration Table Last Updated: Wed May 17 10:06:32 2006

5.17.06



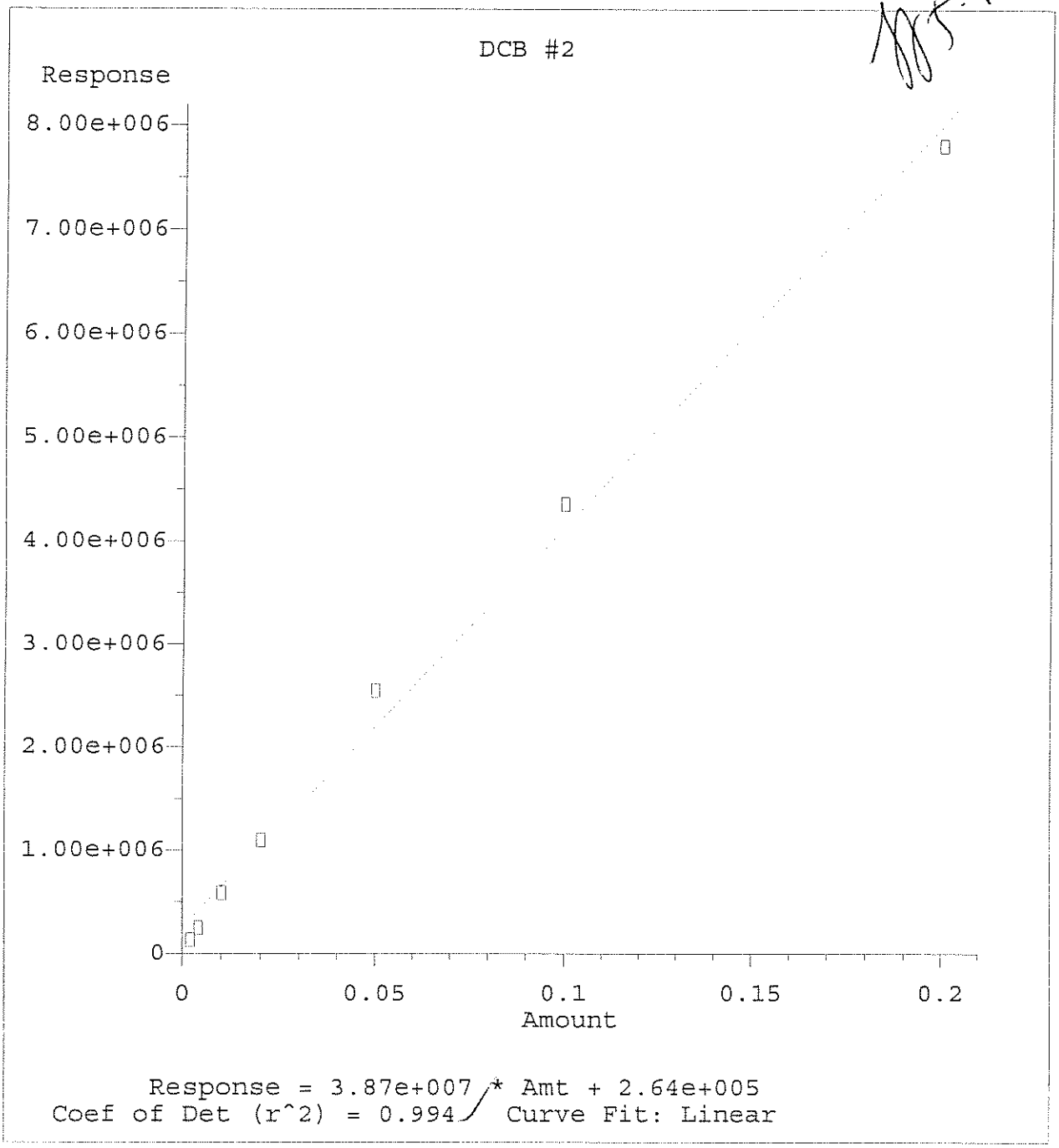
Method Name: C:\HPCHEM\2\METHODS\PBQ05166.M
Calibration Table Last Updated: Wed May 17 10:06:32 2006

5-17-06



Method Name: C:\HPCHEM\2\METHODS\PBQ05166.M
Calibration Table Last Updated: Wed May 17 10:06:32 2006

5-17-06



Method Name: C:\HPCHEM\2\METHODS\PBQ05166.M
Calibration Table Last Updated: Wed May 17 10:06:32 2006

Quantitation Report (Not Reviewed)

Signal #1 : C:\HPCHEM\2\DATA\051606-F\F55868.D\ECD1A.CH Vial: 20
Signal #2 : C:\HPCHEM\2\DATA\051606-F\F55868.D\ECD2B.CH
Acq On : 16 May 2006 9:35 pm Operator:
Sample : P4024 16/60 0.5 CC Inst : Instrumen
Misc : Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 17 10:08 2006 Quant Results File: PBQ05166.RES

Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 10:06:32 2006
Response via : Initial Calibration
DataAcq Meth : PSTQUICK.M

5.17.06 5.17.06

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film

Compound	RT#1	RT#2	Resp#1	Resp#2	ug/ml	ug/ml
System Monitoring Compounds						
1) S TCX	2.64	2.72	3626028	3966484	0.104	0.105
Spiked Amount 0.100	Range 46 - 122		Recovery =		104.00%	105.00%
9) S DCB	10.39	10.04	3115536	4423256	0.104	0.107
Spiked Amount 0.100	Range 26 - 86		Recovery =		104.00%#	107.00%#
Target Compounds						
2) L1 Ar1016 A	3.43	3.37	832389	883397	0.488	0.496
3) L1 Ar1016 B	3.92	3.86	1587624	2003833	0.487	0.500
4) L1 Ar1016 C	4.09	4.02	656707	761021	0.483	0.501
Sum Ar1016 A			3076721	3648252	1.459	1.496
Average Ar1016 A					0.486 97%	0.499 100%
5) L2 Ar1260 A	6.66	6.35	1060994	1236509	0.507	0.507
6) L2 Ar1260 B	7.03	6.82	1233068	1866651	0.493	0.514
7) L2 Ar1260 C	7.92	7.53	954369	1330389	0.482	0.499
8) L2 Ar1260 D	8.49	8.59	934025	2982100	0.481	0.486
Sum Ar1260 A			4182456	7415649	1.963	2.006
Average Ar1260 A					0.491 99%	0.501 100%

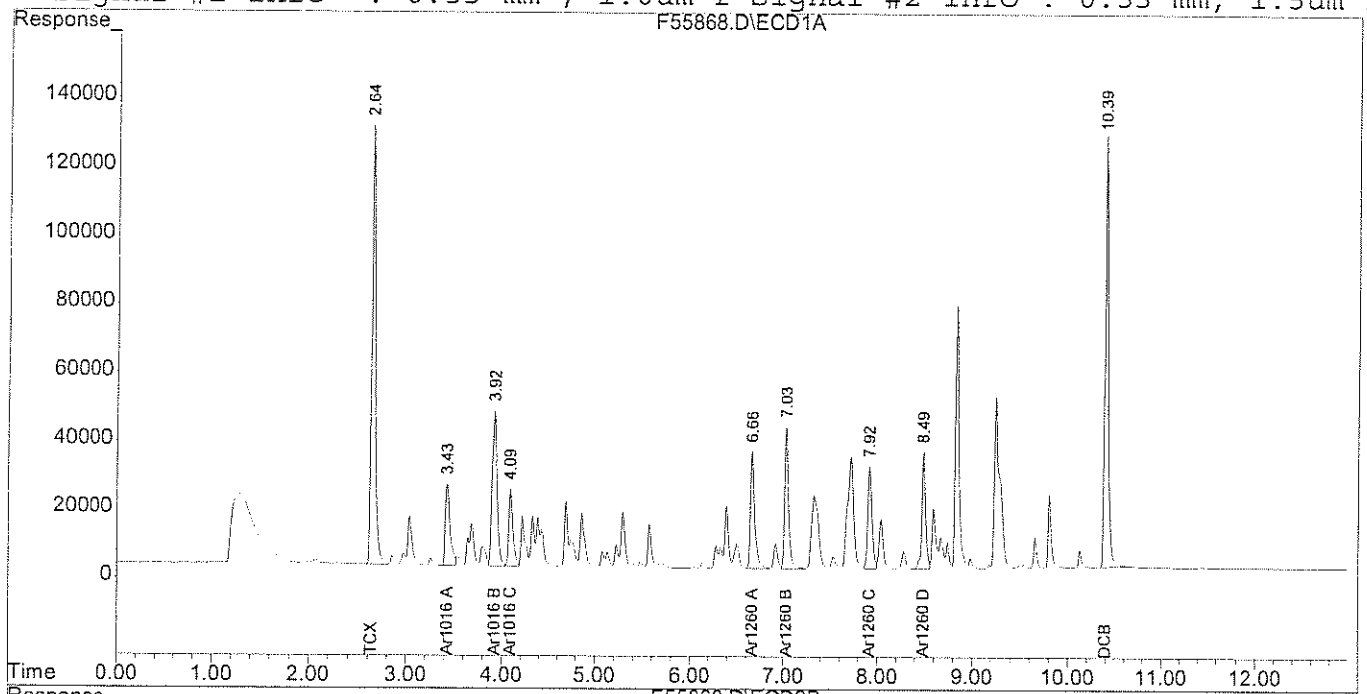
(f)=RT Delta > 1/2 Window (#)=Amounts differ by > 25% (m)=manual int

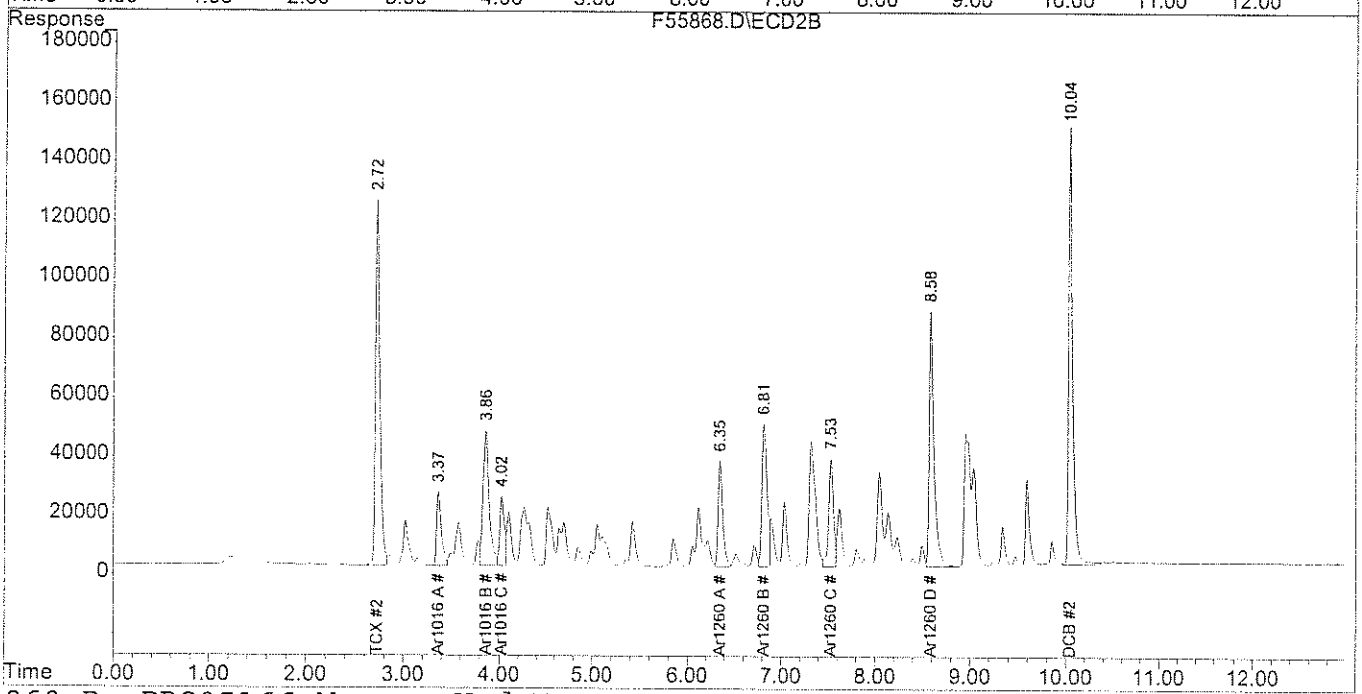
Quantitation Report (Not Reviewed)

Signal #1 : C:\HPCHEM\2\DATA\051606-F\F55868.D\ECD1A.CH Vial: 20
Signal #2 : C:\HPCHEM\2\DATA\051606-F\F55868.D\ECD2B.CH
Acq On : 16 May 2006 9:35 pm Operator:
Sample : P4024 16/60 0.5 CC Inst : Instrumen
Misc : Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 17 10:08 2006 Quant Results File: PBQ05166.RES

Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 10:06:32 2006
Response via : Multiple Level Calibration
DataAcq Meth : PSTQUICK.M

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film





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7A

PCB CONTINUING CALIBRATION CHECK

Site ID: ME

Project No.: GC Column #1: RTX-CLP Pesticides I Data Package ID: 56606

Instrument ID: F GC Column #2: RTX-CLP Pesticides II ID: 0.53 mm

File #: F56057 Cal. Check Date: 5/24/06 15:30

COMPOUND	True Value (ug/mL)	Measured Value (ug/mL)	% Recovery	*	Limits %R	Total Out
PCB 1016	0.25	0.250	100		85-115	0
PCB1260	0.25	0.245	98		85-115	0
PCB 1016#2	0.25	0.266	106		85-115	0
PCB 1260 #2	0.25	0.272	109		85-115	0

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7A

PCB CONTINUING CALIBRATION CHECK

Site ID: ME

Project No.: GC Column #1: RTX-CLP Pesticides I Data Package ID: 56606

Instrument ID: F GC Column #2: RTX-CLP Pesticides II ID: 0.53 mm

File #: F56069 Cal. Check Date: 5/24/06 20:08

COMPOUND	True Value (ug/mL)	Measured Value (ug/mL)	% Recovery	*	Limits %R	Total Out
PCB 1016	0.50	0.492	98		85-115	0
PCB1260	0.50	0.494	99		85-115	0
PCB 1016#2	0.50	0.502	100		85-115	0
PCB 1260 #2	0.50	0.522	104		85-115	0

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7A
PCB CONTINUING CALIBRATION CHECK

Site ID: ME

Project No.: GC Column #1: RTX-CLP Pesticides I Data Package ID: 56606

Instrument ID: F GC Column #2: RTX-CLP Pesticides II ID: 0.53 mm

File #: F56072 Cal. Check Date: 5/25/06 10:32

COMPOUND	True Value (ug/mL)	Measured Value (ug/mL)	% Recovery	*	Limits %R	Total Out
PCB 1016	0.25	0.248	99		85-115	0
PCB1260	0.25	0.228	91		85-115	0
PCB 1016#2	0.25	0.273	109		85-115	0
PCB 1260 #2	0.25	0.288	115		85-115	0

analytics environmental laboratory LLC

7A

PCB CONTINUING CALIBRATION CHECK

Site ID: ME

Project No.: GC Column #1: RTX-CLP Pesticides I Data Package ID: 56606

Instrument ID: F GC Column #2: RTX-CLP Pesticides II ID: 0.53 mm

File #: F56076 Cal. Check Date: 5/25/06 11:45

COMPOUND	True Value (ug/mL)	Measured Value (ug/mL)	% Recovery	*	Limits %R	Total Out
PCB 1016	0.50	0.509	102		85-115	0
PCB1260	0.50	0.482	96		85-115	0
PCB 1016#2	0.50	0.521	104		85-115	0
PCB 1260 #2	0.50	0.551	110		85-115	0

analytics environmental laboratory LLC

PCB SYSTEM MONITORING COMPOUND RETENTION TIME SUMMARY

Quote#: SDG: 56606

Project No: Site ID: Location: ME

GC Columns: RTX-CLPesticidesI ID: 0.32(mm)

Instrument ID: F RTX-CLPesticidesII

Date Analyzed: 5/16/06

Lab File ID (Standard) : F55866 Time Analyzed: 20:51

					COLUMN #1		COLUMN #1	
					SMC1	SMC 2	SMC1	SMC 2
	PESTICIDE STANDARD				2.64	10.39	2.73	10.04
	UPPER LIMIT				2.69	10.44	2.78	10.09
	LOWER LIMIT				2.59	10.34	2.68	9.99
	SAMPLE ID	LAB FILE ID	DATE	TIME	COLUMN #1 SMC1 #	COLUMN #1 SMC 2 #	COLUMN #2 SMC1 #	COLUMN #2 SMC 2 #
01	P4016 16/60 0.01	F55861.D	05/16/06	19:38	2.64	10.39	2.72	10.04
02	P4017 16/60 0.02	F55862.D	05/16/06	19:53	2.64	10.40	2.73	10.04
03	P4018 16/60 0.05	F55863.D	05/16/06	20:07	2.64	10.40	2.73	10.05
04	P4019 16/60 0.1	F55864.D	05/16/06	20:22	2.64	10.40	2.73	10.05
05	P4020 16/60 0.25	F55865.D	05/16/06	20:36	2.64	10.40	2.73	10.04
06	P4021 16/60 0.5	F55866.D	05/16/06	20:51	2.64	10.39	2.73	10.04
07	P4022 16/60 1.0	F55867.D	05/16/06	21:20	2.64	10.40	2.73	10.04
08	P4024 16/60 0.5 CC	F55868.D	05/16/06	21:35	2.64	10.39	2.72	10.04
09	P3883 AR 1221	F55869.D	05/16/06	21:50	NA	NA	NA	NA
10	P3884 AR 1232	F55870.D	05/16/06	22:04	NA	NA	NA	NA
11	P3885 AR 1242	F55871.D	05/16/06	22:19	NA	NA	NA	NA
12	P3886 AR 1248	F55872.D	05/16/06	22:34	NA	NA	NA	NA
13	P3994 AR 1254	F55873.D	05/16/06	22:48	NA	NA	NA	NA
14	P3887 AR 1262	F55874.D	05/16/06	23:03	NA	NA	NA	NA
15	P3888 AR 1268	F55875.D	05/16/06	23:17	NA	NA	NA	NA
16	P4057 AR 1254 0.0	F55876.D	05/16/06	23:47	NA	NA	NA	NA
17	P4056 AR 1254 0.0	F55877.D	05/16/06	0:01	NA	NA	NA	NA
18	P4055 AR 1254 0.1	F55878.D	05/17/06	0:16	NA	NA	NA	NA
19	P4054 AR 1254 0.2	F55879.D	05/17/06	0:31	NA	NA	NA	NA
20	P4053 AR 1254 0.5	F55880.D	05/17/06	0:45	NA	NA	NA	NA
21	P4052 AR 1254 1.0	F55881.D	05/17/06	0:59	NA	NA	NA	NA
22	P4010 AR 1254 0.5	F55882.D	05/17/06	1:13	2.64	10.37	2.72	10.02
23	ZZZZZ	F55883.D	5/17/06	1:28	2.64	10.37	2.72	10.02
24	ZZZZZ	F55885.D	5/17/06	2:12	2.64	10.37	2.72	10.02
25	ZZZZZ	F55886.D	5/17/06	2:26	2.64	10.37	2.72	10.02
26	ZZZZZ	F55895.D	5/17/06	4:53	2.64	10.38	2.73	10.03
27	ZZZZZ	F55896.D	5/17/06	5:07	2.64	10.39	2.73	10.03
28	ZZZZZ	F55897.D	5/17/06	5:22	2.64	10.39	2.73	10.03
29	P4021 16/60 0.5	F55900.D	5/17/06	6:06	2.64	10.39	2.72	10.04
30	P4053 AR 1254 0.5	F55901.D	5/17/06	6:21	NA	NA	NA	NA
31								

analytics environmental laboratory LLC

PCB SYSTEM MONITORING COMPOUND RETENTION TIME SUMMARY

Quote#: SDG: 56606

Project No: Site ID: Location: ME

GC Columns: RTX-CLPesticidesI ID: 0.32(mm)

Instrument ID: F RTX-CLPesticidesII

Date Analyzed: 5/24/06

Lab File ID (Standard) : F56057 Time Analyzed: 15:30

					COLUMN #1		COLUMN #1	
					SMC1	SMC 2	SMC1	SMC 2
	PESTICIDE STANDARD				2.64	10.42	2.72	10.06
	UPPER LIMIT				2.69	10.47	2.77	10.11
	LOWER LIMIT				2.59	10.37	2.67	10.01
	SAMPLE ID	LAB FILE ID	DATE	TIME	COLUMN #1 SMC1 #	COLUMN #1 SMC 2 #	COLUMN #2 SMC1 #	COLUMN #2 SMC 2 #
01	P4020 16/60 0.25	F56057.D	05/24/06	15:30	2.64	10.42	2.72	10.06
02	B05246PAS, A/C	F56058.D	05/24/06	17:13	2.69	10.46	2.77	10.10
03	L05246PASB, A/C	F56059.D	05/24/06	17:28	2.65	10.43	2.73	10.07
04	LD05246PASB, A/C	F56060.D	05/24/06	17:42	2.65	10.42	2.73	10.06
05	56606-1, A/C	F56061.D	05/24/06	17:57	2.64	10.41	2.73	10.06
06	56606-2, A/C	F56062.D	05/24/06	18:11	2.64	10.41	2.72	10.06
07	56606-3, A/C	F56063.D	05/24/06	18:26	2.64	10.40	2.72	10.05
08	56606-4, A/C	F56064.D	05/24/06	18:41	2.64	10.41	2.72	10.05
09	ZZZZZ	F56065.D	05/24/06	19:10	2.64	10.40	2.72	10.05
10	ZZZZZ	F56066.D	05/24/06	19:25	2.64	10.40	2.72	10.05
11	56606-6, MS, A/C	F56067.D	05/24/06	19:39	2.64	10.40	2.72	10.05
12	56606-6, MSD, A/C	F56068.D	05/24/06	19:54	2.63	10.40	2.72	10.05
13	P4021 16/60 0.5	F56069.D	05/24/06	20:08	2.64	10.41	2.72	10.05
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analytics environmental laboratory LLC

PCB SYSTEM MONITORING COMPOUND RETENTION TIME SUMMARY

Quote#: SDG: 56606

Project No: Site ID: Location: ME

GC Columns: RTX-CLPesticidesI ID: 0.32(mm)

Instrument ID: F RTX-CLPesticidesII

Date Analyzed: 5/25/06

Lab File ID (Standard) : F56072 Time Analyzed: 10:32

					COLUMN #1		COLUMN #1	
					SMC1	SMC 2	SMC1	SMC 2
	PESTICIDE STANDARD				2.64	10.43	2.73	10.08
	UPPER LIMIT				2.69	10.48	2.78	10.13
	LOWER LIMIT				2.59	10.38	2.68	10.03
	SAMPLE ID	LAB FILE ID	DATE	TIME	COLUMN #1 SMC1 #	COLUMN #1 SMC 2 #	COLUMN #2 SMC1 #	COLUMN #2 SMC 2 #
01	P4020 16/60 0.25	F56072.D	05/25/06	10:32	2.64	10.43	2.73	10.08
02	B05246PAS,RR A/(	F56073.D	05/25/06	11:01	2.67	10.45	2.75	10.09
03	56606-5, 1:10	F56074.D	05/25/06	11:16	2.64	10.43	2.73	10.07
04	56606-6, 1:2000	F56075.D	05/25/06	11:30	2.64	10.42	2.72	10.06
05	P4021 16/60 0.5	F56076.D	05/25/06	11:45	2.66	10.43	2.73	10.08
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Katahdin ANALYTICAL SERVICES
340 County Road No. 5
P.O. Box 720
Westbrook, ME 04092
Tel: (207) 874-2400
Fax: (207) 775-4029

Analytics

CHAIN of CUSTODY

PLEASE BEAR DOWN AND PRINT LEGIBLY IN PEN

Page ____ of ____

Client: Woodard + Curran	Contact: Karl Kasper	Phone #: (207) 774-2112	Fax #: (207) 774-6635
Address: 41 Hutchins Drive	City: Portland	State: ME	Zip Code: 04102
Purchase Order #: 218736.002	Proj. Name / No.: —		Katahdin Quote #:
Bill (if different than above)	Address		
Sampler (Print / Sign): Josh Bowe		Copies To: Karl Kasper	

LAB USE ONLY

WORK ORDER #:

KATAHDIN PROJECT NUMBER ____

REMARKS: ____

SHIPPING INFO: ☐ FED EX ☐ UPS ☐ CLIENT

AIRBILL NO: ____

TEMP°C 4°C MM ☒ TEMP BLANK ☐ INTACT ☐ NOT INTACT

ANALYSIS AND CONTAINER TYPE PRESERVATIVES

*	Sample Description	Date / Time coll'd	Matrix	No. of Cntrs.	Filt. ☐Y ☐N	Filt. ☐Y ☐N	Filt. ☐Y ☐N	Filt. ☐Y ☐N	Filt. ☐Y ☐N	Filt. ☐Y ☐N	Filt. ☐Y ☐N	Filt. ☐Y ☐N	Filt. ☐Y ☐N	Filt. ☐Y ☐N
					Unpreserved 4 oz									
	SS01	5/23/06 / 15:50	Soil	1	PCB					5606		-1		
	SS02	5/23/06 / 15:05	Soil	↓	↓							-2		
	SS03	5/23/06 / 15:25	Soil	↓	↓							-3		
	SS04	/ 15:35	Soil	↓	↓							-4		
	SS13	/ 16:00	Soil	↓	↓							-5		
	SS14	/ 16:03	Soil	↓	↓							-6		
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COMMENTS: See Karl for sample Turn-around time ✓ 05/24/06

Relinquished By: (Signature)	Date / Time	Received By: (Signature)	Relinquished By: (Signature)	Date / Time	Received By: (Signature)
Josh Bowe	5/23 19:13				
Wendy Faenzer	5/24 10:49 am	M. Martisens			

95/95

THE TERMS AND CONDITIONS ON THE REVERSE SIDE HEREOF SHALL GOVERN

analytics environmental laboratory LLC

195 Commerce Way Suite E
Portsmouth, New Hampshire 03801
603-436-5111 Fax 603-430-2151
800-929-9906
www.analyticslab.com

June 6, 2006

Mr. Karl Kasper
Woodard & Curran, Inc.
41 Hutchins drive
Portland ME 04102

RE: **Analytical Results Case Narrative**
Analytics # 56617

Dear Mr. Kasper;

Enclosed please find the analytical results for samples submitted for the above-mentioned project. The attached Cover Page lists the sample IDs, Lab tracking numbers and collection dates for the samples included in this deliverable.

Samples were analyzed Polychlorinated Biphenyls (PCBs) by EPA Method 8082.

Unless otherwise noted in the Non-conformance Summary listed below, all of the quality control (QC) criteria including initial calibration, calibration verification, surrogate recovery, holding time and method accuracy/precision for these analyses were within acceptable limits.

This Level III data package has been assembled in the following order:

Case Narrative/Non-Conformance Summary
Sample Log Sheet - Cover Page
PCB Form 1 Data Sheet for Samples and Blanks
PCB Form 10 Confirmation Results Form (if required)
PCB Form 2 Surrogate Recoveries
PCB Form 3 MS/MSD (LCS) Recoveries
PCB Form 4 Method Blank Summary (equiv.) GC/MS Logbook Sheets
PCB Form 6 Initial Calibration Data
PCB Form 7 Continuing Calibration Check
PCB Form 8 System Monitoring Compound RT Summary
Chain of Custody (COC) Forms

QC NON CONFORMANCE SUMMARY

Sample Receipt:
No QC deviations.

PCBs by EPA Method 8082:
No extra sample was provided for MS/MSD analysis for the wipe matrix. Laboratory control samples were used to assess accuracy and precision.

Sample 56606-6 (SS14) was selected for MS/MSD analysis however the data wasn't usuable due to the high concentrations of PCB 1260 detected in the sample. The laboratory control samples for the batch (L05246PASB/LD05246PASB) were in control for all analytes and results were reported without qualification.

Sample 56617-5 (SB-11 (0-4) was reported with a PCB 1016 and PCB 1260 detected. Due to interferences in the sample the laboratory was unable to determine whether the PCB in the sample was PCB 1016 or PCB 1242. The laboratory quantitated the result and reported as PCB 1016 with an "*" qualifier.

Sample 617-32 (W-8) had % difference between the analytical columns above the laboratory acceptance criteria. Results were reported with a "D" flag and a comment to this affect.

The continuing calibration standard (File# F56132) analyzed 05/26/06 had PCB 1260 recovery (116%) above the acceptance criteria of 115% on column #2. PCB 1260 had acceptable recovery on column #1. Results for samples associated with this standard were reported off column#1.

Several samples had retention times for one of the surrogates slightly outside the established windows. Results were reported without qualification.

Sincerely,
ANALYTICS Environmental Laboratory, LLC

Stephen Knollmeyer
Laboratory Director

analytics environmental laboratory LLC

195 Commerce Way Suite E
Portsmouth, New Hampshire 03801
603-436-5111 Fax 603-430-2151
800-929-9906
www.analyticslab.com

Mr. Karl Kasper
Woodard & Curran
41 Hutchins Drive
Portland ME 04102

Report Number: 56617

Revision: Rev. 0

Re: GP **218736**

Enclosed are the results of the analyses on your sample(s). Samples were received on 25 May 2006 and analyzed for the tests listed below. Samples were received in acceptable condition, with the exceptions noted below or on the chain of custody. The results reported herein conform to the most current NELAC standards, where applicable, unless otherwise narrated in the body of the report. Please see individual reports for specific methodologies and references.

Lab Number	Sample Date	Station Location	Analysis	Comments
56617-1	05/24/06	SB-2 (0-4)	EPA 8082 (PCBs only)	
56617-2	05/24/06	SB-3 (0-4)	EPA 8082 (PCBs only)	
56617-3	05/24/06	SB-4 (0-4)	EPA 8082 (PCBs only)	
56617-4	05/24/06	SB-10 (0-1)	EPA 8082 (PCBs only)	
56617-5	05/24/06	SB-11 (0-4)	EPA 8082 (PCBs only)	
56617-6	05/24/06	SB-12 (0-4)	EPA 8082 (PCBs only)	
56617-7	05/24/06	SB-15 (0-4)	EPA 8082 (PCBs only)	
56617-8	05/24/06	SB-18 (0-4)	EPA 8082 (PCBs only)	
56617-9	05/24/06	SB-25 (0-4)	EPA 8082 (PCBs only)	
56617-10	05/24/06	SB-26 (0-4)	EPA 8082 (PCBs only)	
56617-11	05/24/06	SB-28 (0-4)	EPA 8082 (PCBs only)	
56617-12	05/24/06	SC-01	EPA 8082 (PCBs only)	
56617-13	05/24/06	SC-02	EPA 8082 (PCBs only)	
56617-14	05/24/06	SC-03	EPA 8082 (PCBs only)	
56617-15	05/24/06	SS-05	EPA 8082 (PCBs only)	

Sample Receipt Exceptions: None

Analytics Environmental Laboratory is certified by the states of New Hampshire, Maine, Massachusetts, Connecticut, Rhode Island, North Carolina, New York, Virginia, Pennsylvania and is validated by the U.S. Army Corps of Engineers (MRD) and U.S. Navy (NFESC). A list of actual certified parameters is available upon request.

If you have any further question on the analytical methods or these results, do not hesitate to call.

Authorized signature _____
Stephen L. Knollmeyer Lab. Director

Date 05/26/06

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analytics environmental laboratory LLC

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Lab Number	Sample Date	Station Location	Analysis	Comments
56617-16	05/24/06	SS-06	EPA 8082 (PCBs only)	
56617-17	05/24/06	SS-07	EPA 8082 (PCBs only)	
56617-18	05/24/06	SS-08	EPA 8082 (PCBs only)	
56617-19	05/24/06	SS-09	EPA 8082 (PCBs only)	
56617-20	05/24/06	SS-09 DUP	EPA 8082 (PCBs only)	
56617-21	05/24/06	SS-10	EPA 8082 (PCBs only)	
56617-22	05/24/06	SS-11	EPA 8082 (PCBs only)	
56617-23	05/24/06	SS-12	EPA 8082 (PCBs only)	
56617-24	05/24/06	W-1	EPA 8082 (PCB's only)	
56617-25	05/24/06	W-2	EPA 8082 (PCB's only)	
56617-26	05/24/06	W-3	EPA 8082 (PCB's only)	
56617-27	05/24/06	W-4	EPA 8082 (PCB's only)	
56617-28	05/24/06	W-5	EPA 8082 (PCB's only)	
56617-29	05/24/06	W-6	EPA 8082 (PCB's only)	
56617-30	05/24/06	W-6 DUP	EPA 8082 (PCB's only)	

Sample Receipt Exceptions: None

Analytics Environmental Laboratory is certified by the states of New Hampshire, Maine, Massachusetts, Connecticut, Rhode Island, North Carolina, New York, Virginia, Pennsylvania and is validated by the U.S. Army Corps of Engineers (MRD) and U.S. Navy (NFESC). A list of actual certified parameters is available upon request.

If you have any further question on the analytical methods or these results, do not hesitate to call.

Authorized signature Melissa Hall for
Stephen L. Knollmeyer Lab. Director

Date 05/26/06

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analytics environmental laboratory LLC

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Mr. Karl Kasper
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Report Number: 56617

Revision: Rev. 0

Re: GP

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Lab Number	Sample Date	Station Location	Analysis	Comments
56617-31	05/24/06	W-7	EPA 8082 (PCB's only)	
56617-32	05/24/06	W-8	EPA 8082 (PCB's only)	
56617-33	05/24/06	W-9	EPA 8082 (PCB's only)	
56617-34	05/24/06	W-10	EPA 8082 (PCB's only)	
56617-35	05/24/06	W-11	EPA 8082 (PCB's only)	
56617-36	05/24/06	W-12	EPA 8082 (PCB's only)	
56617-37	05/24/06	W-13	EPA 8082 (PCB's only)	

Sample Receipt Exceptions: None

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Authorized signature _____
Stephen L. Knollmeyer Lab. Director

Date 05/26/06

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analytics environmental laboratory LLC

EXTRACTION
LOGBOOK PAGES

Transfer Date: 05/25/06

Initials: SC

EXTRACT TRANSFER SHEET

Organochlorine PST

Sample	Analysis	Extraction Date	Matrix	Initial Volume(mL)	Initial Weight(g)	Final Volume(mL)	% Solids	Extraction Comments
56617-1	8082ASE	05/25/06	Solid		15.54	10.0		
56617-2	8082ASE	05/25/06	Solid		15.98	10.0		
56617-3	8082ASE	05/25/06	Solid		15.85	10.0		
56617-4	8082ASE	05/25/06	Solid		15.63	10.0		
56617-5	8082ASE	05/25/06	Solid		15.27	10.0		
56617-6	8082ASE	05/25/06	Solid		15.31	10.0		
56617-7	8082ASE	05/25/06	Solid		15.94	10.0		
56617-8	8082ASE	05/25/06	Solid		15.65	10.0		
56617-9	8082ASE	05/25/06	Solid		15.26	10.0		
56617-10	8082ASE	05/25/06	Solid		15.21	10.0		
56617-11	8082ASE	05/25/06	Solid		15.21	10.0		
56617-12	8082ASE	05/25/06	Solid		15.06	10.0		
56617-13	8082ASE	05/25/06	Solid		15.46	10.0		
56617-14	8082ASE	05/25/06	Solid		4.43	10.0		

VBM 5/25/06

AK 5.25.06

Pesticide and PCB Solid and Waste Preparation Logbook

Method References SW-846 3550B Sonication
SW-846 3545 Accelerated Solvent Extraction (ASE)
SW-846 3580 Concentrated Waste Dilution

Date: 5-25-06 Time: 1:00pm QC Batch ID : Q-P-052406-1-SC
Initials: SL Prep. Batch ID P-P-052406-1-SC
Reviewed by: JM 5/25/06

Sample #	Matrix	Weight (g)	mL Surr	mL Spike	Final Vol Hexane	Cook Down Date	Vessel #	CleanUp
B052506 PAS	Na2SO4	15.00	1.0	—	10mL	5/25/06	30403	ASE
L057526 PASB	Si Sand	15.19	1.0	1.0		SL	38773	
56617-1	Soil	15.54	1.0	—			38336	
-2		15.98	1.0	—			30439	
-3		15.85	1.0	—			30812	
-4		15.63	1.0	—			30937	
-5		15.27	1.0	—			31141-104	
-6		15.31	1.0	—			30454	
-7		15.94	1.0	—			31144-112	
-8		15.65	1.0	—			38746	
-9		15.26	1.0	—			39730	
-10		15.21	1.0	—			31179	
-11		15.21	1.0	—			31180	
-12	Concrete	15.06	1.0	—			38837	
-13		15.46	1.0	—			31159	
-14		14.43	1.0	—			38879	
		SL 5/25/06						
							5/25/06 SL	

Hexane Lot Number: 40060
Si Sand Lot Number: 074697 (1)
Na_2SO_4 Lot Number: 45251602
1:1 hexane and Acetone /
Diatomaceous earth /
Conc H2SO4 B41044
Florosil Cartridge

PP Surrogate (1.0 ug/mL):	Lot # P4012	Exp Date 11.10.06
(2) PCB MS (4.0 ug/mL):	P3992	9.21.06
Pest Spike (500 ug/L):		

Transfer Date: 05/26/06

Initials: SC

EXTRACT TRANSFER SHEET

Organochlorine PST

Sample	Analysis	Extraction Date	Matrix	Initial Volume(mL)	Initial Weight(g)	Final Volume(mL)	% Solids	Extraction Comments
56617-15	8082ASE	05/26/06	Solid		15.97	10.0	82.73	
56617-16	8082ASE	05/26/06	Solid		15.61	10.0	92.60	
56617-17	8082ASE	05/26/06	Solid		15.28	10.0	95.37	
56617-18	8082ASE	05/26/06	Solid		15.15	10.0	93.92	
56617-19	8082ASE	05/26/06	Solid		15.53	10.0	94.24	
56617-20	8082ASE	05/26/06	Solid		15.37	10.0	89.74	
56617-21	8082ASE	05/26/06	Solid		15.28	10.0	78.97	
56617-22	8082ASE	05/26/06	Solid		15.74	10.0	88.29	
56617-23	8082ASE	05/26/06	Solid		15.13	10.0	89.68	

5.26.06

AK 5.26.06

Pesticide and PCB Solid and Waste Preparation Logbook

Method References SW-846 3550B Sonication
SW-846 3545 Accelerated Solvent Extraction (ASE)
SW-846 3580 Concentrated Waste Dilution

Date: 5-26-06 Time: 6:00AM QC Batch ID : Q-P-052606-1SL
Initials: SL Prep. Batch ID D-P-052606-1-SL
Reviewed by: AK 5.26.06

Sample #	Matrix	Weight (g)	mL Surr (1)	mL Spike (2)/(3)	Final Vol Hexane	Cook Down Date	Vessel #	CleanUp
B052666PAS	Na_2SO_4	15.00g	1.0	–	10mL	5/26/06	39241	ALL
LU52666PASB	Si Sand	15.02g	1.0	1.0 (2)		SL	39286	
LD052666PASB	↓	15.96g	1.0	1.0 (2)			38925	
56617-15	Soil	15.97g	1.0	–			31212	
-16		15.61g	1.0	–			38720	
-17		15.28g	1.0	–	10mL	5/26/06	39163	
-18		15.15g	1.0	–		AK	38901	
-19		15.53g	1.0	–			30787	
-20		15.37g	1.0	–			38845	
-21		15.28g	1.0	–			39385	
-22		15.74g	1.0	–			38680	
-23		15.13g	1.0	–			31203	
-23, MS		15.02g	1.0	1.0 (2)			38930	
-23, MSD	↓	13.05g*	1.0	1.0 (2)			30531	

5/26/06 SL

Hexane Lot Number: 46060
Si Sand Lot Number: 074617 (1)
Na_2SO_4 Lot Number: 45251602 (2)
1:1 hexane and Acetone /
Diatomaceous earth /
Conc H2SO4 B41041
Florosil Cartridge

	Lot #	Exp Date
PP Surrogate (1.0 ug/mL):	P4012	11.D06
PCB MS (4.0 ug/mL):	P3992	9.21.06
Pest Spike (500 ug/L):	/	/

* used up all remaining sample

03/18/05 Rev 1

Transfer Date: 05/25/06

Initials: RM

EXTRACT TRANSFER SHEET

Organochlorine PST

Sample	Analysis	Extraction Date	Matrix	Initial Volume(mL)	Initial Weight(g)	Final Volume(mL)	% Solids	Extraction Comments
56617-24	PCBWIPE	05/25/06	Wipe		1.00	25.0 ✓		
56617-25	PCBWIPE	05/25/06	Wipe		1.00	25.0 ✓		
56617-26	PCBWIPE	05/25/06	Wipe		1.00	25.0 ✓		
56617-27	PCBWIPE	05/25/06	Wipe		1.00	25.0 ✓		
56617-28	PCBWIPE	05/25/06	Wipe		1.00	25.0 ✓		
56617-29	PCBWIPE	05/25/06	Wipe		1.00	25.0 ✓		
56617-30	PCBWIPE	05/25/06	Wipe		1.00	25.0 ✓		
56617-31	PCBWIPE	05/25/06	Wipe		1.00	25.0 ✓		
56617-32	PCBWIPE	05/25/06	Wipe		1.00	25.0 ✓		
56617-33	PCBWIPE	05/25/06	Wipe		1.00	25.0 ✓		
56617-34	PCBWIPE	05/25/06	Wipe		1.00	25.0 ✓		
56617-35	PCBWIPE	05/25/06	Wipe		1.00	25.0 ✓		
56617-36	PCBWIPE	05/25/06	Wipe		1.00	25.0 ✓		
56617-37	PCBWIPE	05/25/06	Wipe		1.00	25.0 ✓		

RM 5/25/06

✓ AK 5.25.06

Pesticide and PCB Solid and Waste Preparation Logbook

PCB WIPE

Method References SW-846 3550B Sonication
SW-846 3545 Accelerated Solvent Extraction (ASE)
SW-846 3580 Concentrated Waste Dilution

Date: 5/25/06 Time: 2:00 QC Batch ID : Q-P-052506-1-RM
Initials: RM Prep. Batch ID P-P-052506-1-RM
Reviewed by: AK 5.25.06

Sample #	Matrix	Weight (g)	mL (1) Surr	mL (2) Spike	Final Vol Hexane	Cook Down Date	Vessel #	CleanUp
B052506PWIPE	wipe	—	1.0	—	25mL	5/25/06	N/A	A/C
L052506PWIPE		—	1.0	1.0		SC		
LD052506PWIPE		—	1.0	1.0				
56617-24		—	1.0	—				
-25		—	1.0	—				
-26		—	1.0	—				
-27		—	1.0	—				
-28		—	1.0	—				
-29		—	1.0	—				
-30		—	1.0	—				
-31		—	1.0	—				
-32		—	1.0	—				
-33		—	1.0	—				
-34		—	1.0	—				
-35		—	1.0	—				
-36		—	1.0	—				
-37		—	1.0	—				

5/25/06 SC

Hexane Lot Number: 46060 (1)
Si Sand Lot Number: —
Na_2SO_4 Lot Number: —
1:1 hexane and Acetone —
Diatomaceous earth —
Conc H2SO4 B411041
Florosil Cartridge —

	Lot #	Exp Date
PP Surrogate (1.0 ug/mL):	P4012	11-10-06
(2) PCB MS (4.0 ug/mL):	P3992	9-21-06
Pest Spike (500 ug/L):	—	—

analytics environmental laboratory LLC

PCB
DATA SUMMARIES

analytics environmental laboratory LLC

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Portland ME 04102

June 6, 2006

CLIENT SAMPLE ID

Project Name: GP

Project Number: 218736

Field Sample ID: LABQC

SAMPLE DATA

Lab Sample ID: B05246PAS
Matrix: Solid
Percent Solid: 100
Dilution Factor: 1.0
Collection Date:
Lab Receipt Date:
Extraction Date: 05/24/06
Analysis Date: 05/24/06

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit µg/kg	Results µg/kg
PCB-1016	15	U
PCB-1221	15	U
PCB-1232	15	U
PCB-1242	15	U
PCB-1248	15	U
PCB-1254	15	U
PCB-1260	15	U

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	84	%
Decachlorobiphenyl	68	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample was analyzed according to "Test Methods for Evaluating Solid Waste, SW-846 Method 8081."

COMMENTS: Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3545. Results are expressed on a dry weight basis.

PCB Report

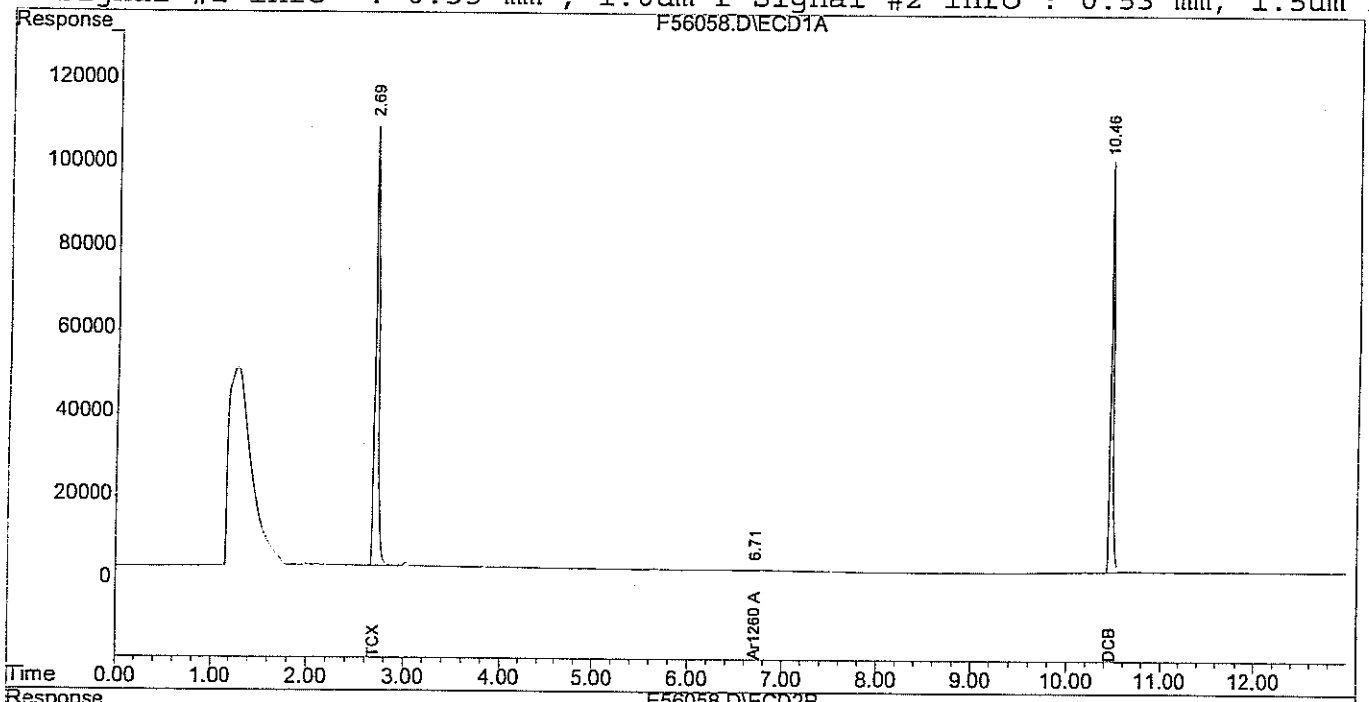
Authorized signature

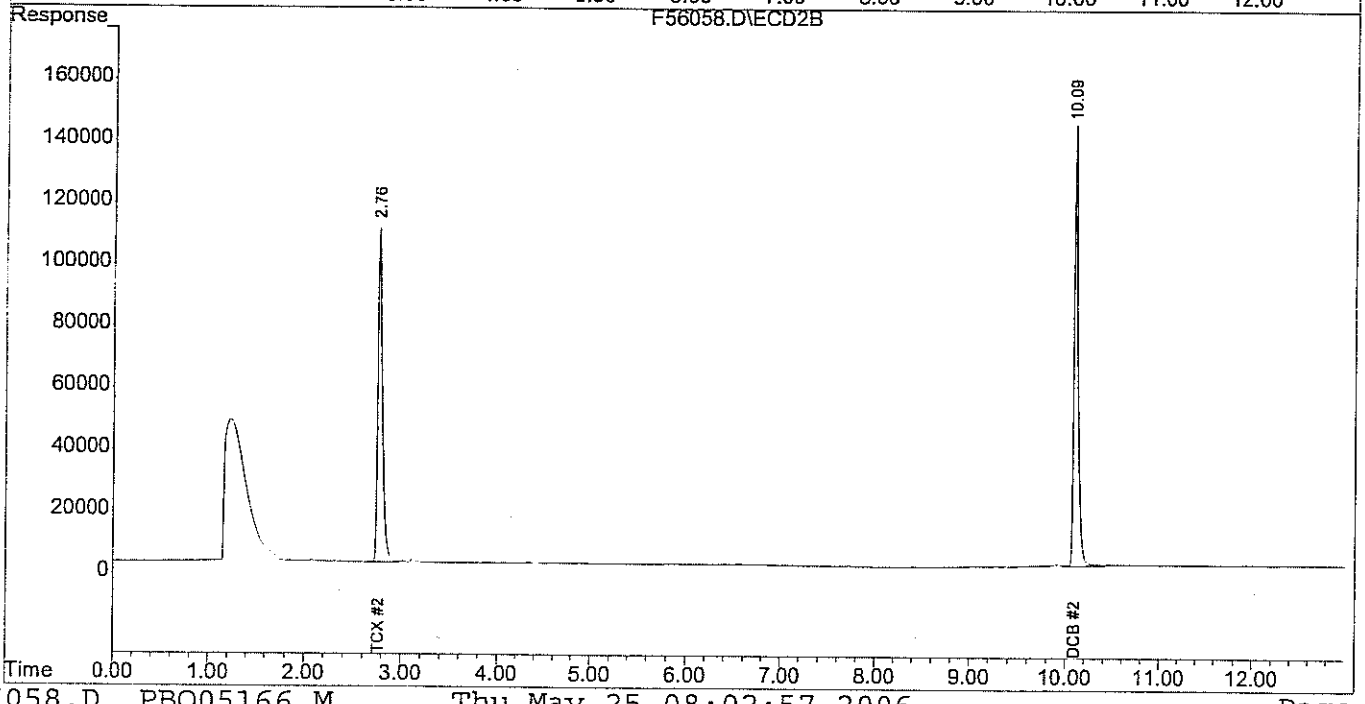
Quantitation Report (Not Reviewed)

Signal #1 : C:\HPCHEM\2\DATA\052406-F\F56058.D\ECD1A.CH Vial: 2
Signal #2 : C:\HPCHEM\2\DATA\052406-F\F56058.D\ECD2B.CH
Acq On : 24 May 2006 5:13 pm Operator:
Sample : B05246PAS, A/C Inst : Instrumen
Misc : SOIL Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 25 8:02 2006 Quant Results File: PBQ05166.RES

Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 10:06:32 2006
Response via : Multiple Level Calibration
DataAcq Meth : PSTQUICK.M

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film





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Portland ME 04102

June 6, 2006

CLIENT SAMPLE ID

Project Name: GP

Project Number: 218736

Field Sample ID: W-1

SAMPLE DATA

Lab Sample ID: B0525PWIPE
Matrix: Wipe
Percent Solid: N/A
Dilution Factor: 1.0
Collection Date:
Lab Receipt Date:
Extraction Date: 05/25/06
Analysis Date: 05/25/06

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit µg/wipe	Results µg/wipe
PCB-1016	0.5	U
PCB-1221	0.5	U
PCB-1232	0.5	U
PCB-1242	0.5	U
PCB-1248	0.5	U
PCB-1254	0.5	U
PCB-1260	0.5	U

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	93	%
Decachlorobiphenyl	100	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8081/82.

Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3550B.

COMMENTS:

PCB Report

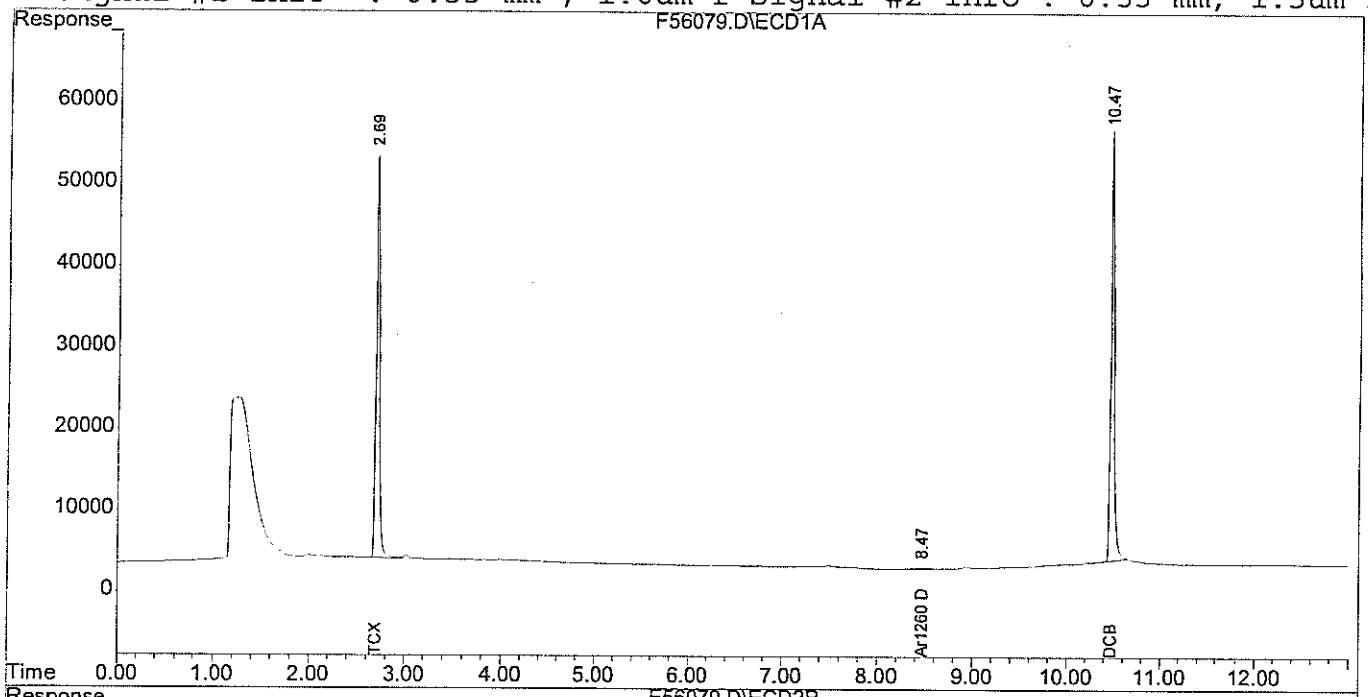
Authorized signature

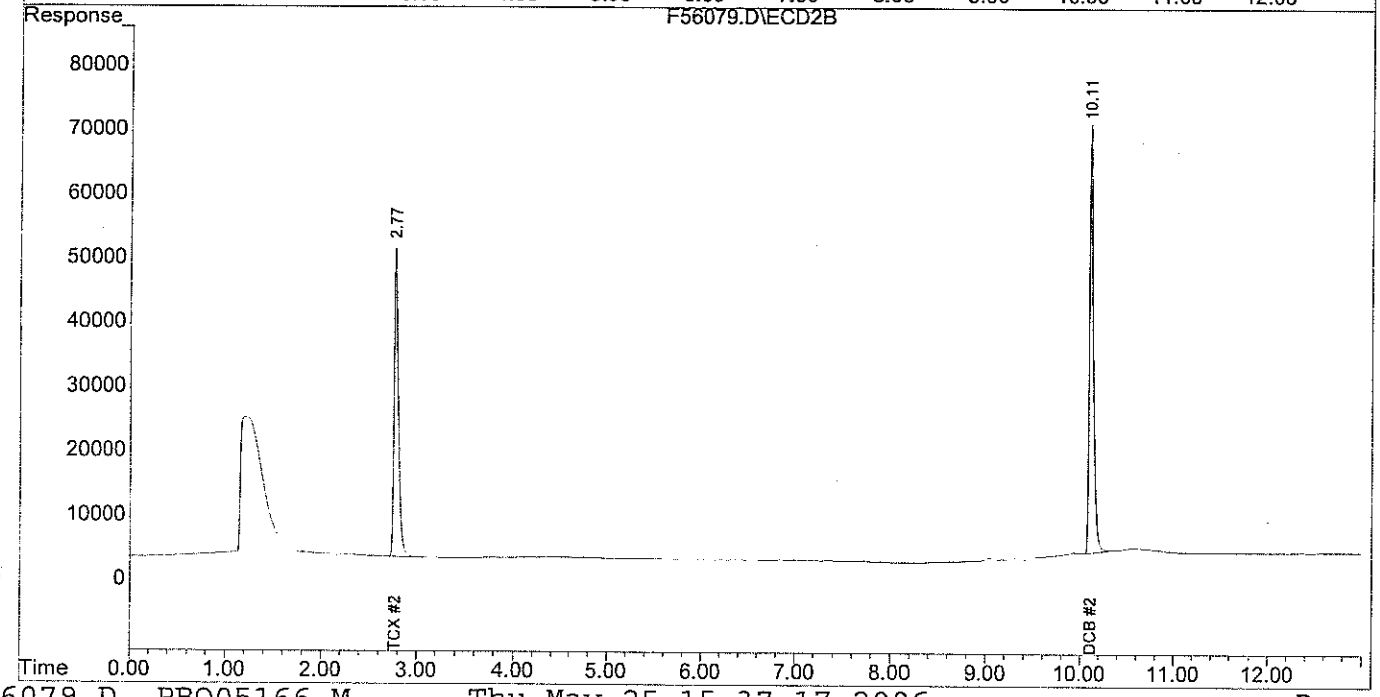
Quantitation Report (QT Reviewed)

Signal #1 : C:\HPCHEM\2\DATA\052506-F\F56079.D\ECD1A.CH Vial: 2
Signal #2 : C:\HPCHEM\2\DATA\052506-F\F56079.D\ECD2B.CH
Acq On : 25 May 2006 3:10 pm Operator:
Sample : B05256PWIPE, A/C Inst : Instrumen
Misc : SOIL Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 25 15:37 2006 Quant Results File: PBQ05166.RES

Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 10:06:32 2006
Response via : Multiple Level Calibration
DataAcq Meth : PSTQUICK.M

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film





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Portland ME 04102

June 6, 2006

SAMPLE DATA

CLIENT SAMPLE ID

Project Name: GP

Project Number: 218736

Field Sample ID: LABQC

Lab Sample ID: B05256PAS
Matrix: Solid
Percent Solid: 100
Dilution Factor: 1.0
Collection Date:
Lab Receipt Date:
Extraction Date: 05/25/06
Analysis Date: 05/25/06

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit µg/kg	Results µg/kg
PCB-1016	15	U
PCB-1221	15	U
PCB-1232	15	U
PCB-1242	15	U
PCB-1248	15	U
PCB-1254	15	U
PCB-1260	15	U

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	99	%
Decachlorobiphenyl	109	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample was analyzed according to "Test Methods for Evaluating Solid Waste, SW-846 Method 8081."

COMMENTS: Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3545. Results are expressed on a dry weight basis.

PCB Report

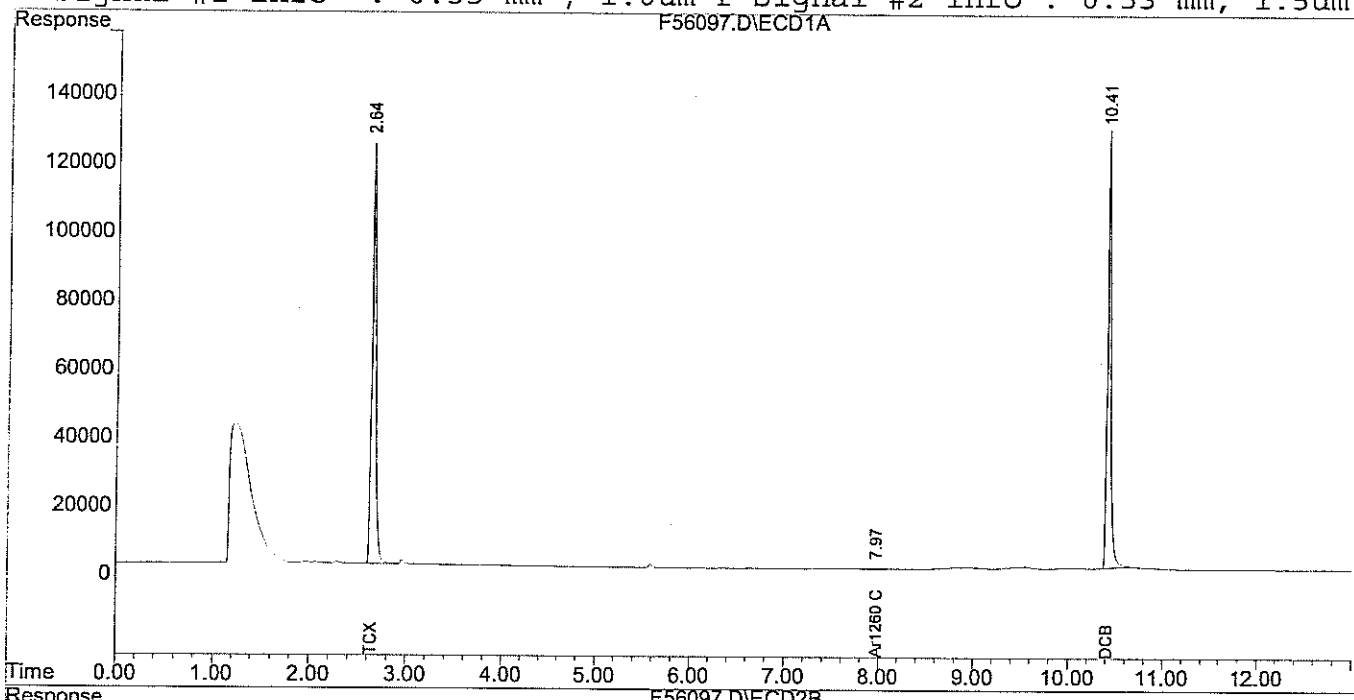
Authorized signature

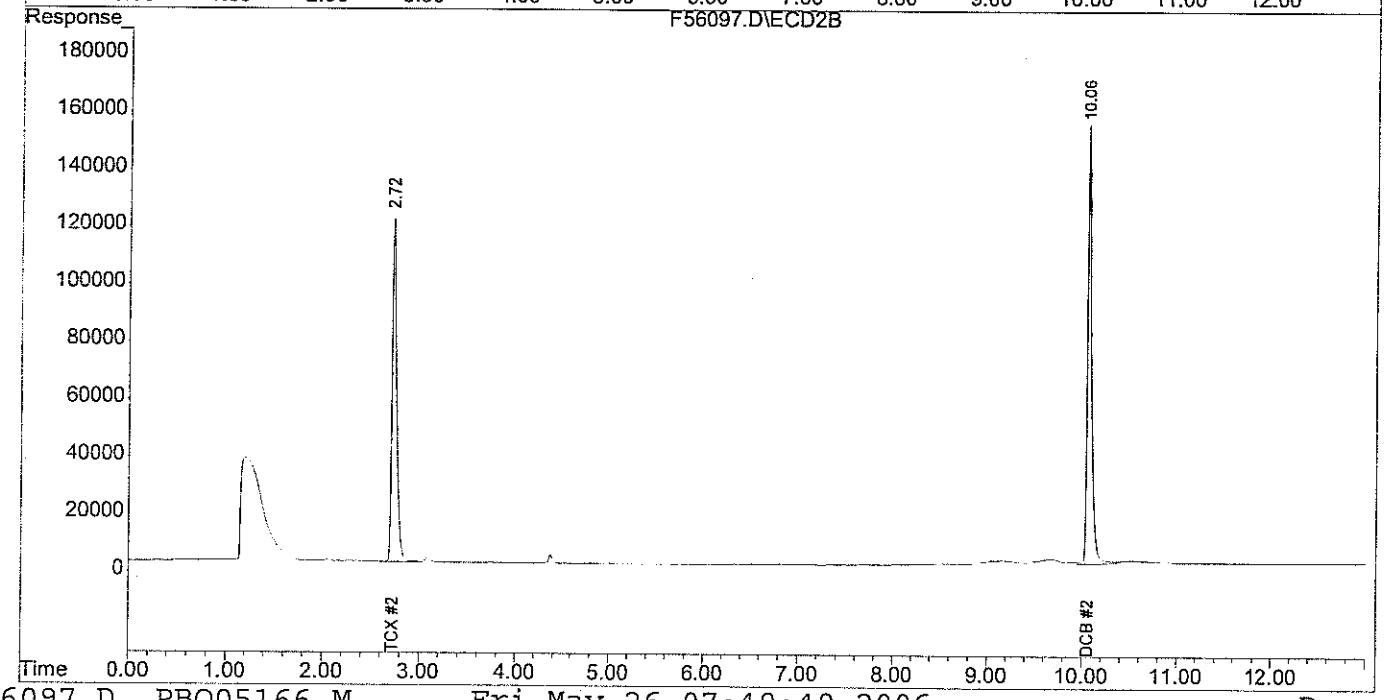
Quantitation Report (Not Reviewed)

Signal #1 : C:\HPCHEM\2\DATA\052506-F\F56097.D\ECD1A.CH Vial: 20
Signal #2 : C:\HPCHEM\2\DATA\052506-F\F56097.D\ECD2B.CH
Acq On : 25 May 2006 8:17 pm Operator:
Sample : B05256PAS, A/C Inst : Instrumen
Misc : SOIL Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 26 7:49 2006 Quant Results File: PBQ05166.RES

Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 10:06:32 2006
Response via : Multiple Level Calibration
DataAcq Meth : PSTQUICK.M

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film





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Portland ME 04102

June 6, 2006

SAMPLE DATA

CLIENT SAMPLE ID

Project Name: GP

Project Number: 218736

Field Sample ID: LABQC

Lab Sample ID: B05266PAS
Matrix: Solid
Percent Solid: 100
Dilution Factor: 1.0
Collection Date:
Lab Receipt Date:
Extraction Date: 05/26/06
Analysis Date: 05/26/06

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit µg/kg	Results µg/kg
PCB-1016	15	U
PCB-1221	15	U
PCB-1232	15	U
PCB-1242	15	U
PCB-1248	15	U
PCB-1254	15	U
PCB-1260	15	U

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	102	%
Decachlorobiphenyl	99	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample was analyzed according to "Test Methods for Evaluating Solid Waste, SW-846 Method 8081."

COMMENTS: Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3545. Results are expressed on a dry weight basis.

PCB Report

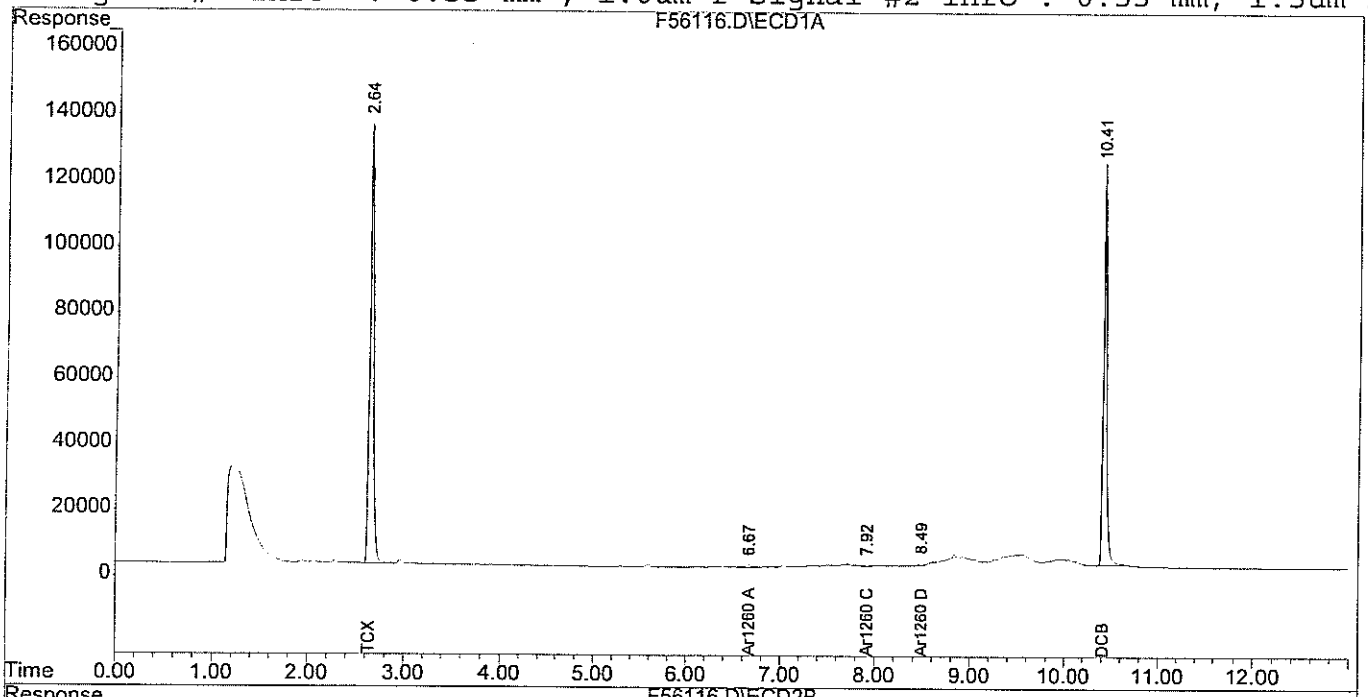
Authorized signature Melena Sali

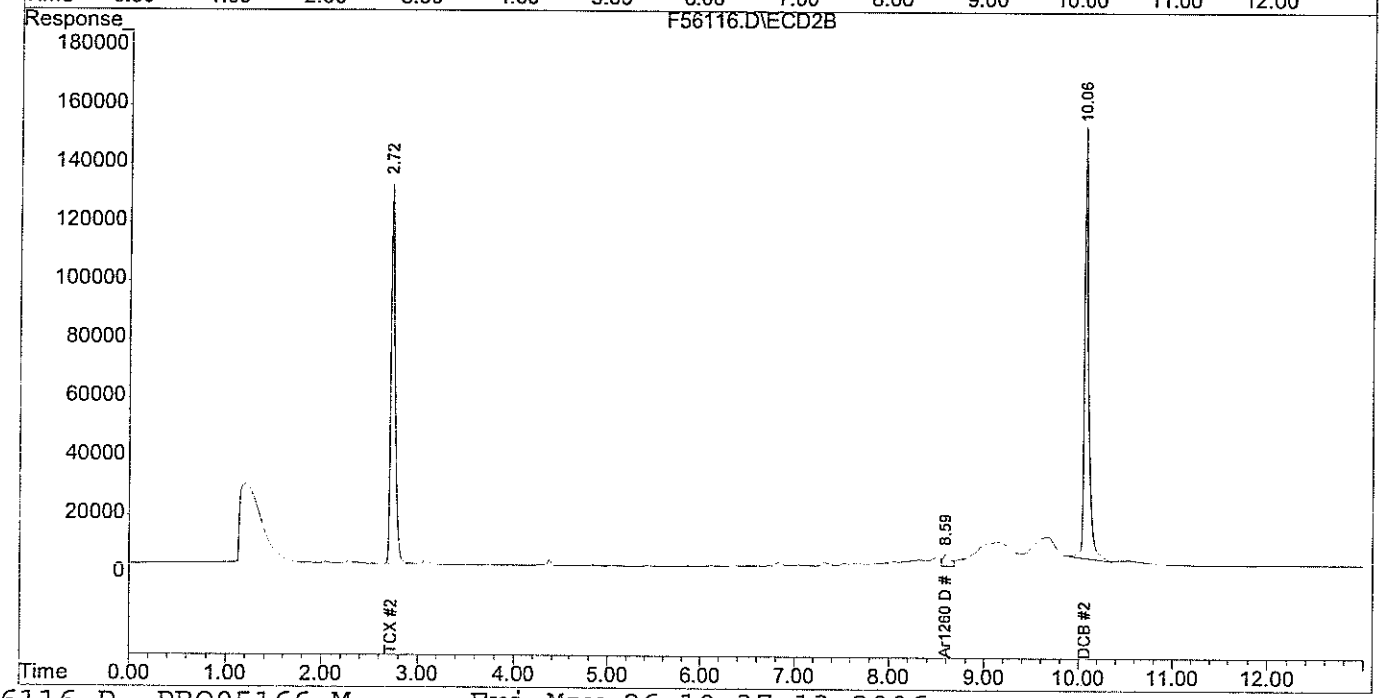
Quantitation Report (QT Reviewed)

Signal #1 : C:\HPCHEM\2\DATA\052506-F\F56116.D\ECD1A.CH Vial: 36
Signal #2 : C:\HPCHEM\2\DATA\052506-F\F56116.D\ECD2B.CH
Acq On : 26 May 2006 7:30 am Operator:
Sample : B05266PAS, A/C IRR 56600 Inst : Instrumen
Misc : SOIL Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 26 10:37 2006 Quant Results File: PBQ05166.RES

Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 10:06:32 2006
Response via : Multiple Level Calibration
DataAcq Meth : PSTQUICK.M

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film





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June 6, 2006

CLIENT SAMPLE ID

Project Name: GP

Project Number: 218736

Field Sample ID: LABQC

SAMPLE DATA

Lab Sample ID: B05266PAS RR
Matrix: Solid
Percent Solid: 100
Dilution Factor: 1.0
Collection Date:
Lab Receipt Date:
Extraction Date: 05/26/06
Analysis Date: 05/26/06

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit µg/kg	Results µg/kg
PCB-1016	15	U
PCB-1221	15	U
PCB-1232	15	U
PCB-1242	15	U
PCB-1248	15	U
PCB-1254	15	U
PCB-1260	15	U

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	109	%
Decachlorobiphenyl	112	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample was analyzed according to "Test Methods for Evaluating Solid Waste, SW-846 Method 8081."

COMMENTS: Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3545. Results are expressed on a dry weight basis.

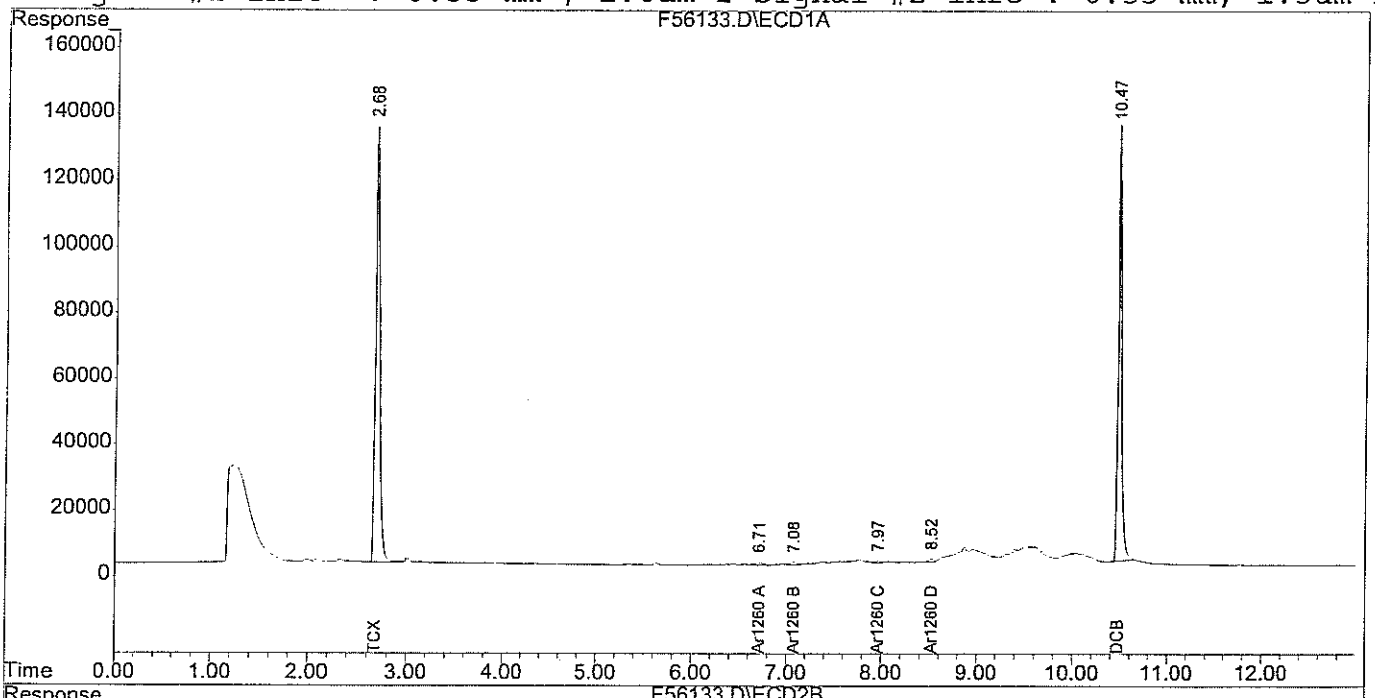
PCB Report

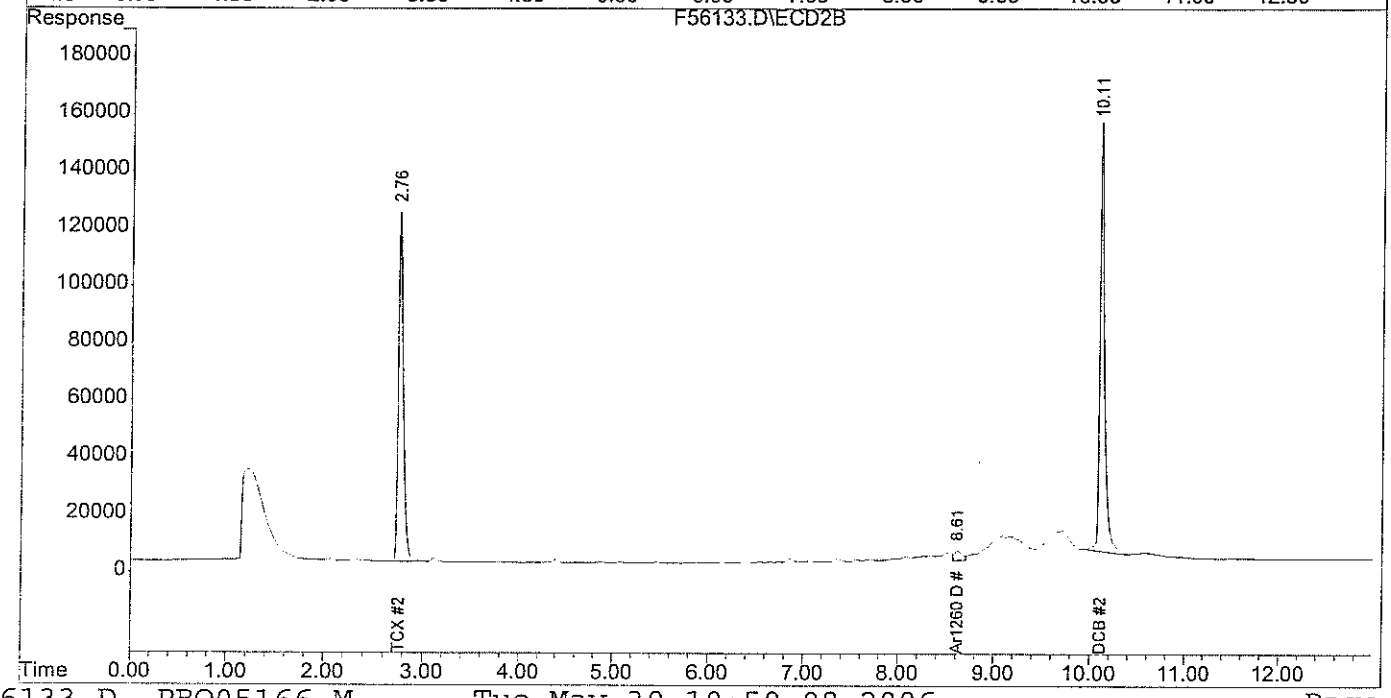
Authorized signature

Signal #1 : C:\HPCHEM\2\DATA\052606-F\F56133.D\ECD1A.CH Vial: 36
Signal #2 : C:\HPCHEM\2\DATA\052606-F\F56133.D\ECD2B.CH
Acq On : 26 May 2006 2:07 pm Operator:
Sample : B05266PAS, A/C, RR Inst : Instrume
Misc : SOIL Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 30 10:50 2006 Quant Results File: PBQ05166.RES

Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 10:06:32 2006
Response via : Multiple Level Calibration
DataAcq Meth : PSTQUICK.M

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film





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May 26, 2006

CLIENT SAMPLE ID

Project Name: GP

Project Number: 218736

Field Sample ID: SB-2 (0-4)

SAMPLE DATA

Lab Sample ID: 56617-1
Matrix: Solid
Percent Solid: 81
Dilution Factor: 1.2
Collection Date: 05/24/06
Lab Receipt Date: 05/25/06
Extraction Date: 05/25/06
Analysis Date: 05/25/06

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit µg/kg	Results µg/kg
PCB-1016	18	U
PCB-1221	18	U
PCB-1232	18	U
PCB-1242	18	U
PCB-1248	18	U
PCB-1254	18	U
PCB-1260	18	U

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	91	%
Decachlorobiphenyl	96	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082.

Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3545.

COMMENTS: Results are expressed on a dry weight basis.

PCB Report

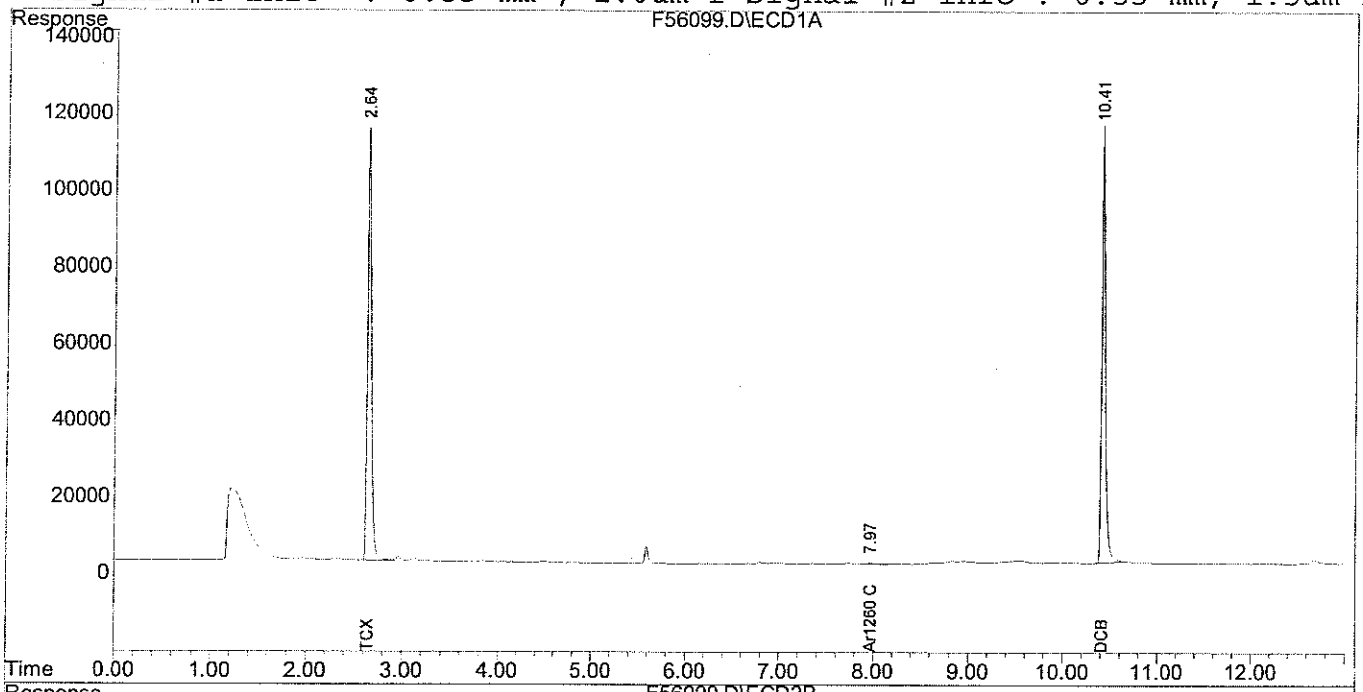
Authorized signature Milena Tull

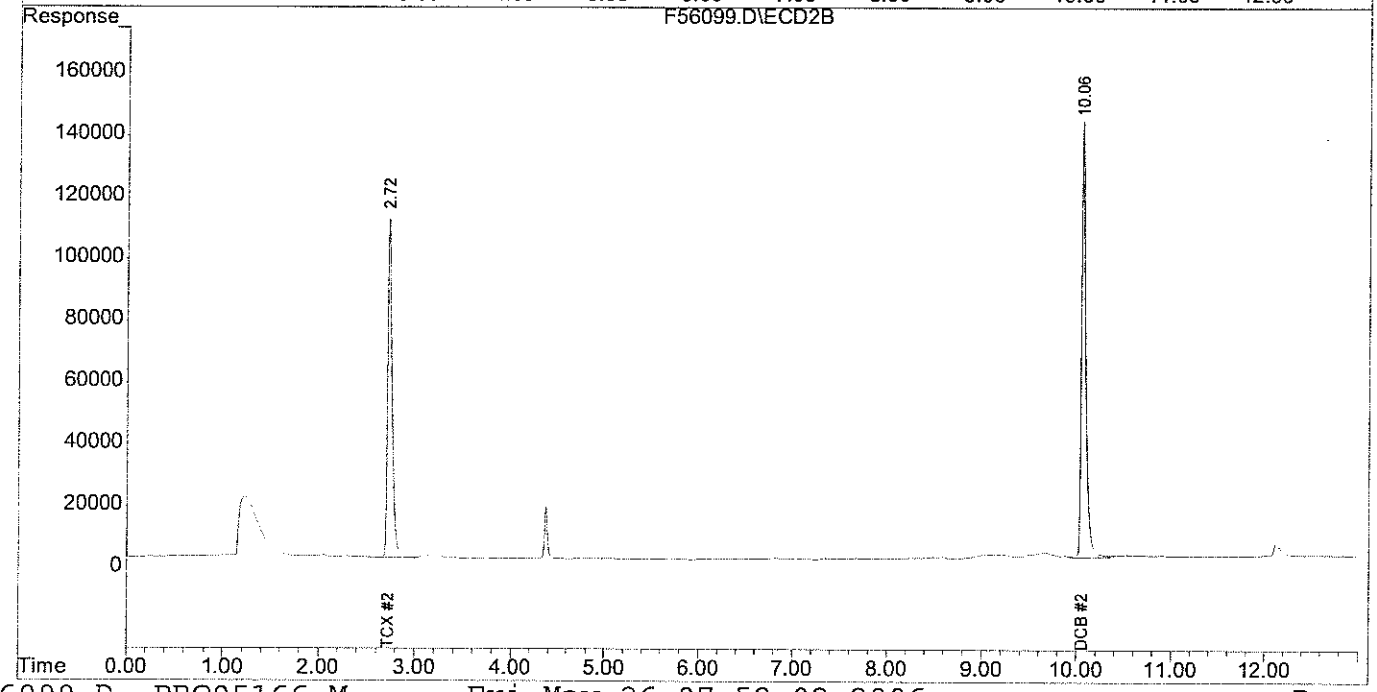
Quantitation Report (QT Reviewed)

Signal #1 : C:\HPCHEM\2\DATA\052506-F\F56099.D\ECD1A.CH Vial: 22
Signal #2 : C:\HPCHEM\2\DATA\052506-F\F56099.D\ECD2B.CH
Acq On : 25 May 2006 8:46 pm Operator:
Sample : 56617-1, A/C Inst : Instrumen
Misc : SOIL Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 26 7:51 2006 Quant Results File: PBQ05166.RES

Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 10:06:32 2006
Response via : Multiple Level Calibration
DataAcq Meth : PSTQUICK.M

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film





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Mr. Karl Kasper
Woodard & Curran
41 Hutchins Drive
Portland ME 04102

May 26, 2006

CLIENT SAMPLE ID

Project Name: GP

Project Number: 218736

Field Sample ID: SB-3 (0-4)

SAMPLE DATA

Lab Sample ID: 56617-2
Matrix: Solid
Percent Solid: 91
Dilution Factor: 1.0
Collection Date: 05/24/06
Lab Receipt Date: 05/25/06
Extraction Date: 05/25/06
Analysis Date: 05/25/06

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit µg/kg	Results µg/kg
PCB-1016	15	U
PCB-1221	15	U
PCB-1232	15	U
PCB-1242	15	U
PCB-1248	15	U
PCB-1254	15	**276**
PCB-1260	15	**282**

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	91	%
Decachlorobiphenyl	86	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082.

Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3545.

COMMENTS: Results are expressed on a dry weight basis.

PCB Report

Authorized signature Mylena Mull

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PCB CONFIRMATION RESULTS

SDG: 56617

Project No: Site ID: Location: ME

Instrument ID: F GC Columns: RTX-CLPesticidesI ID: 0.32(mm)
RTX-CLPesticidesII

Sample ID: 56617-2

COMPOUND	RESPONSE #1 ug/kg	RESPONSE #2 ug/kg	DIFFERENCE %	#	% DIFFERENCE Acceptance Limit
PCB 1254	276	294	6		40
PCB 1260	282	330	16		40

Column to be used to flag % Difference with an asterisk

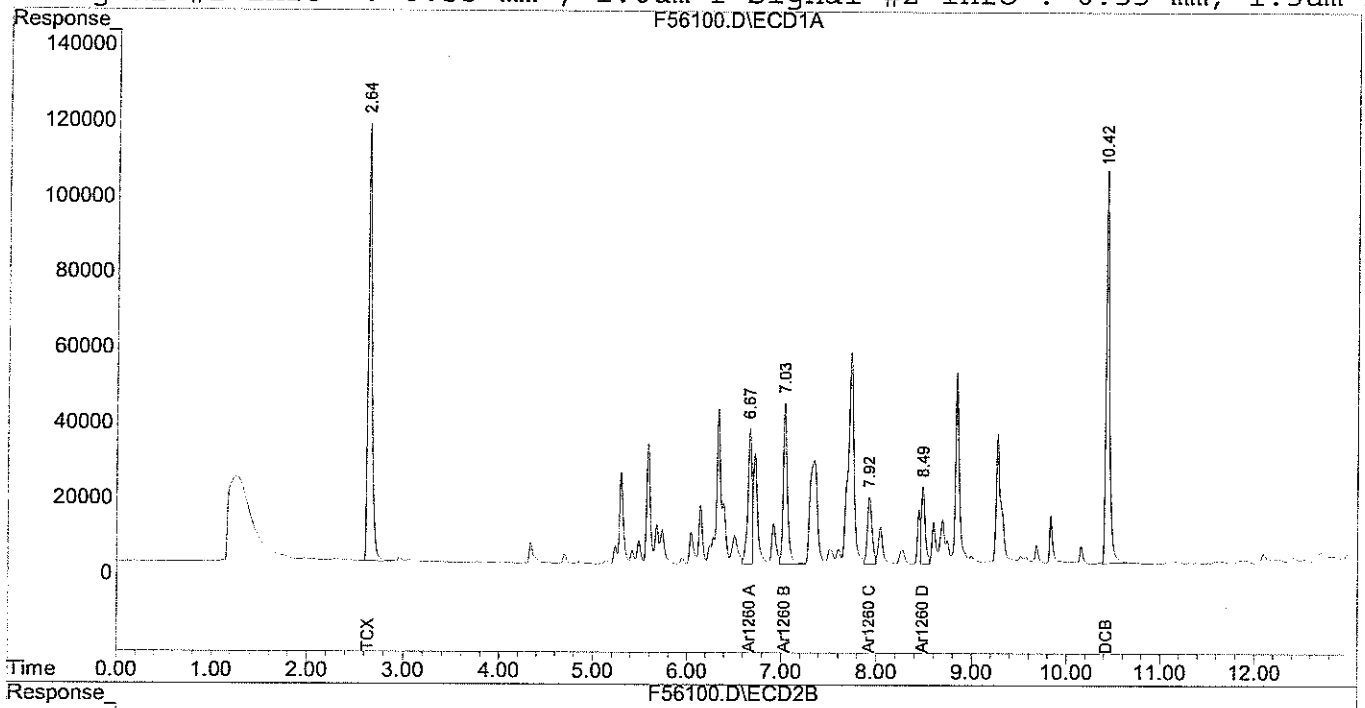
* Values outside of QC limits

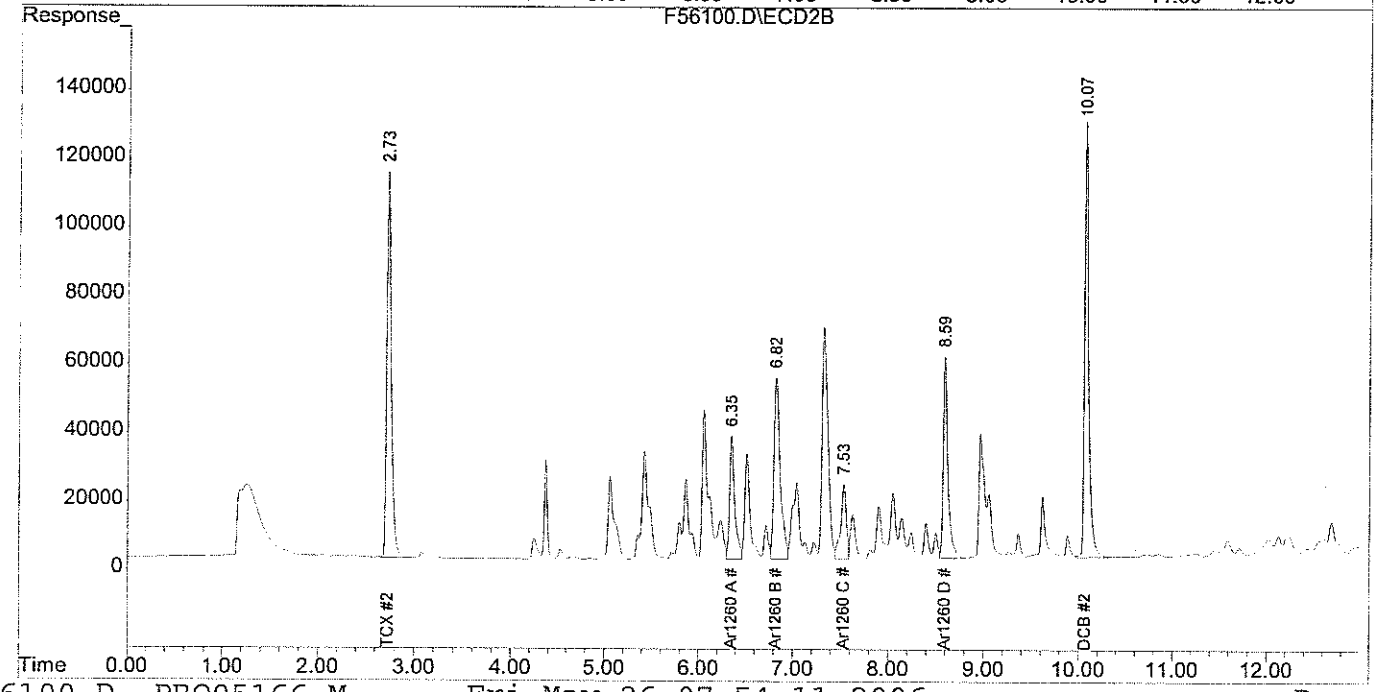
Quantitation Report (QT Reviewed)

Signal #1 : C:\HPCHEM\2\DATA\052506-F\F56100.D\ECD1A.CH Vial: 23
Signal #2 : C:\HPCHEM\2\DATA\052506-F\F56100.D\ECD2B.CH
Acq On : 25 May 2006 9:01 pm Operator:
Sample : 56617-2, A/C Inst : Instrumen
Misc : SOIL Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 26 7:53 2006 Quant Results File: PBQ05166.RES

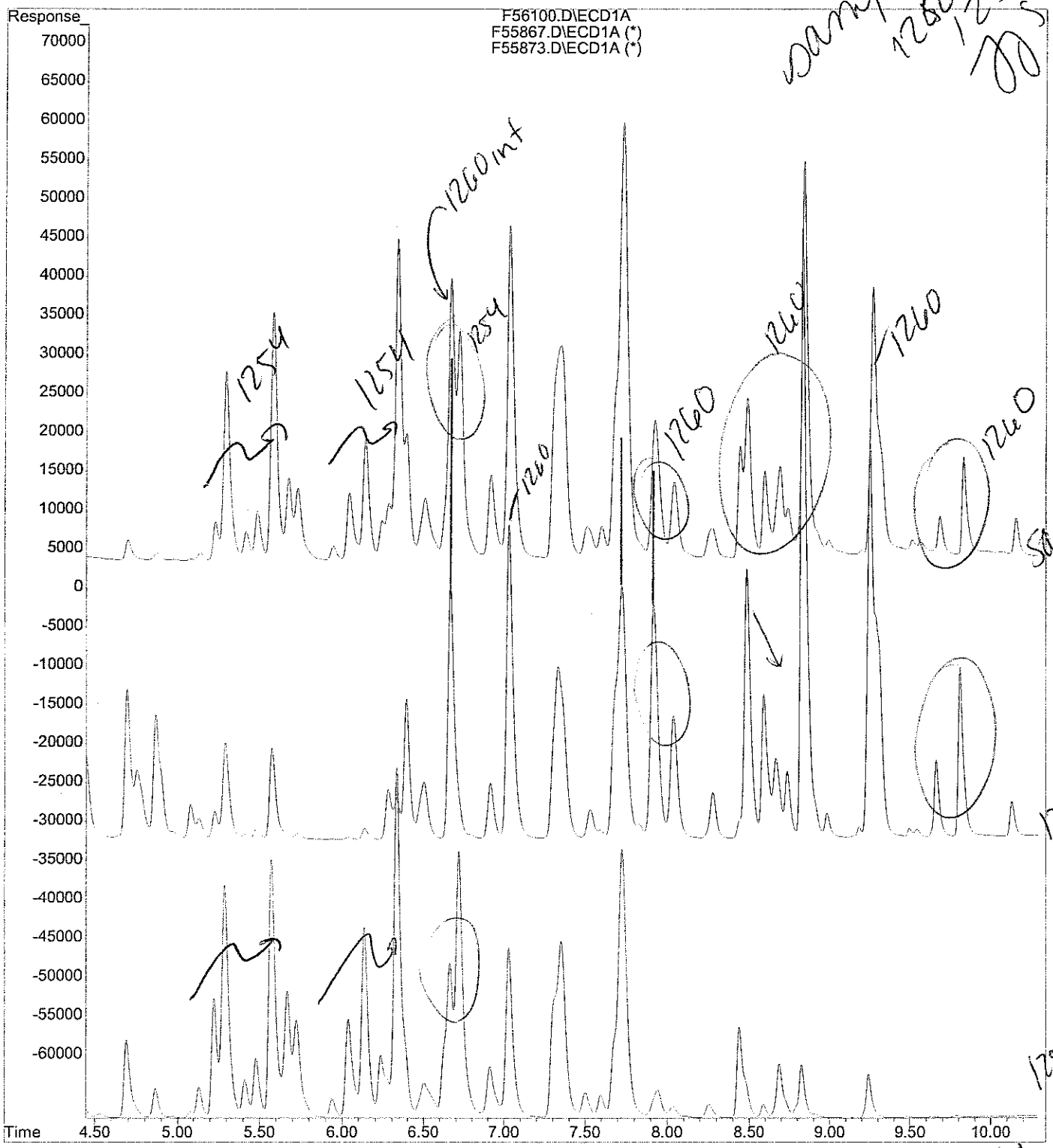
Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 10:06:32 2006
Response via : Multiple Level Calibration
DataAcq Meth : PSTQUICK.M

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film





File : C:\HPCHEM\2\DATA\052506-F\F56100.D
Operator :
Acquired : 25 May 2006 9:01 pm using AcqMethod PSTQUICK.M
Instrument : Instrumen
Sample Name: 56617-2, A/C
Misc Info : SOIL
Vial Number: 23



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May 30, 2006

CLIENT SAMPLE ID

Project Name: GP

Project Number: 218736

Field Sample ID: SB-4 (0-4)

SAMPLE DATA

Lab Sample ID: 56617-3 DL
Matrix: Solid
Percent Solid: 93
Dilution Factor: 5
Collection Date: 05/24/06
Lab Receipt Date: 05/25/06
Extraction Date: 05/25/06
Analysis Date: 05/26/06

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit µg/kg	Results µg/kg
PCB-1016	75	U
PCB-1221	75	U
PCB-1232	75	U
PCB-1242	75	U
PCB-1248	75	U
PCB-1254	75	**200**
PCB-1260	75	**306**

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	90	%
Decachlorobiphenyl	90	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082.

Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3545.

COMMENTS: Results are expressed on a dry weight basis.

PCB Report

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PCB CONFIRMATION RESULTS

SDG: 56617

Project No: Site ID: Location: ME

Instrument ID: F GC Columns: RTX-CLPesticidesI ID: 0.32(mm)
RTX-CLPesticidesII

Sample ID: 56617-3

COMPOUND	RESPONSE #1 ug/kg	RESPONSE #2 ug/kg	DIFFERENCE %	#	% DIFFERENCE Acceptance Limit
PCB 1254	200	177	13		40
PCB 1260	306	323	5		40

Column to be used to flag % Difference with an asterisk

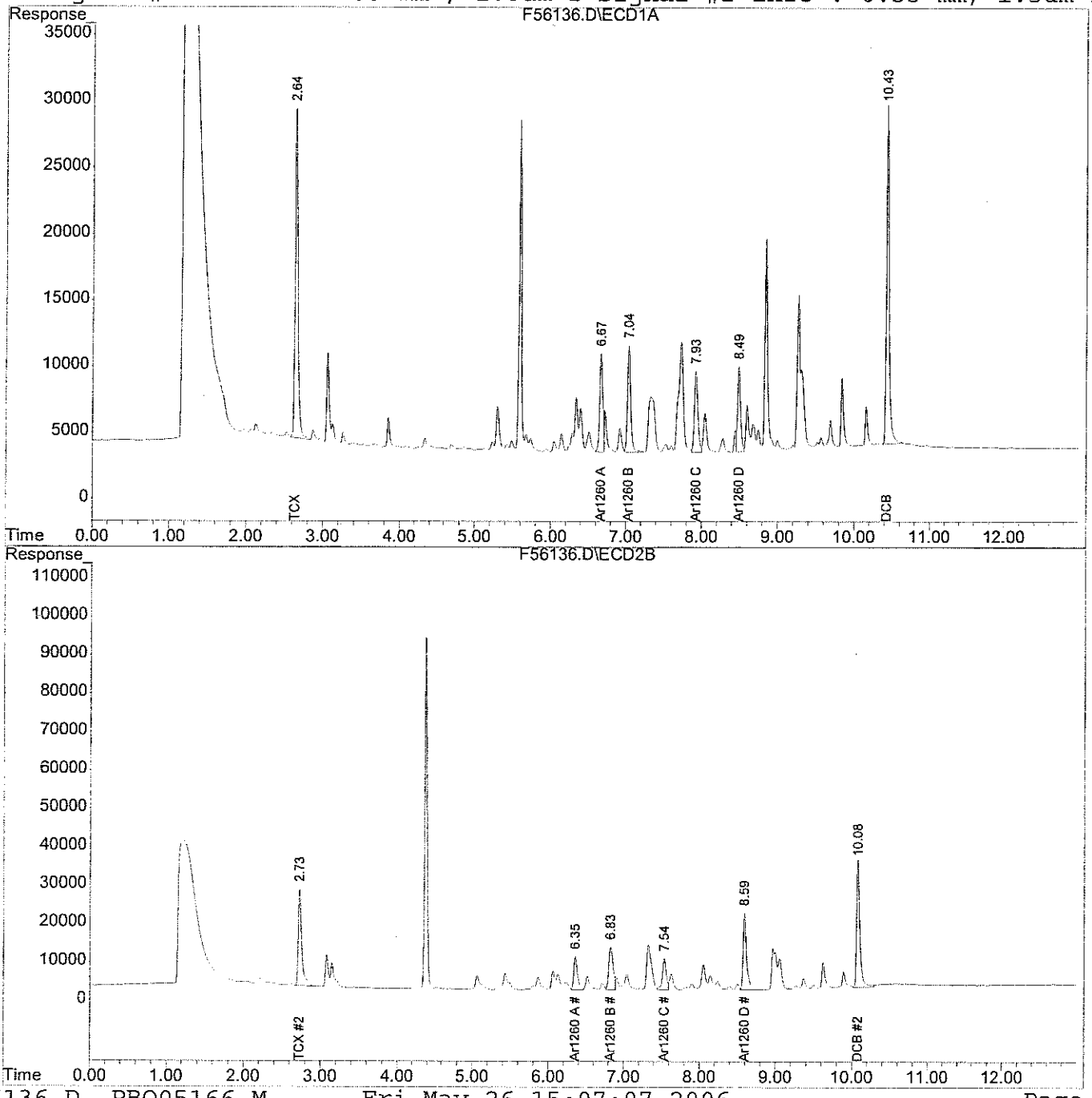
* Values outside of QC limits

Quantitation Report (QT Reviewed)

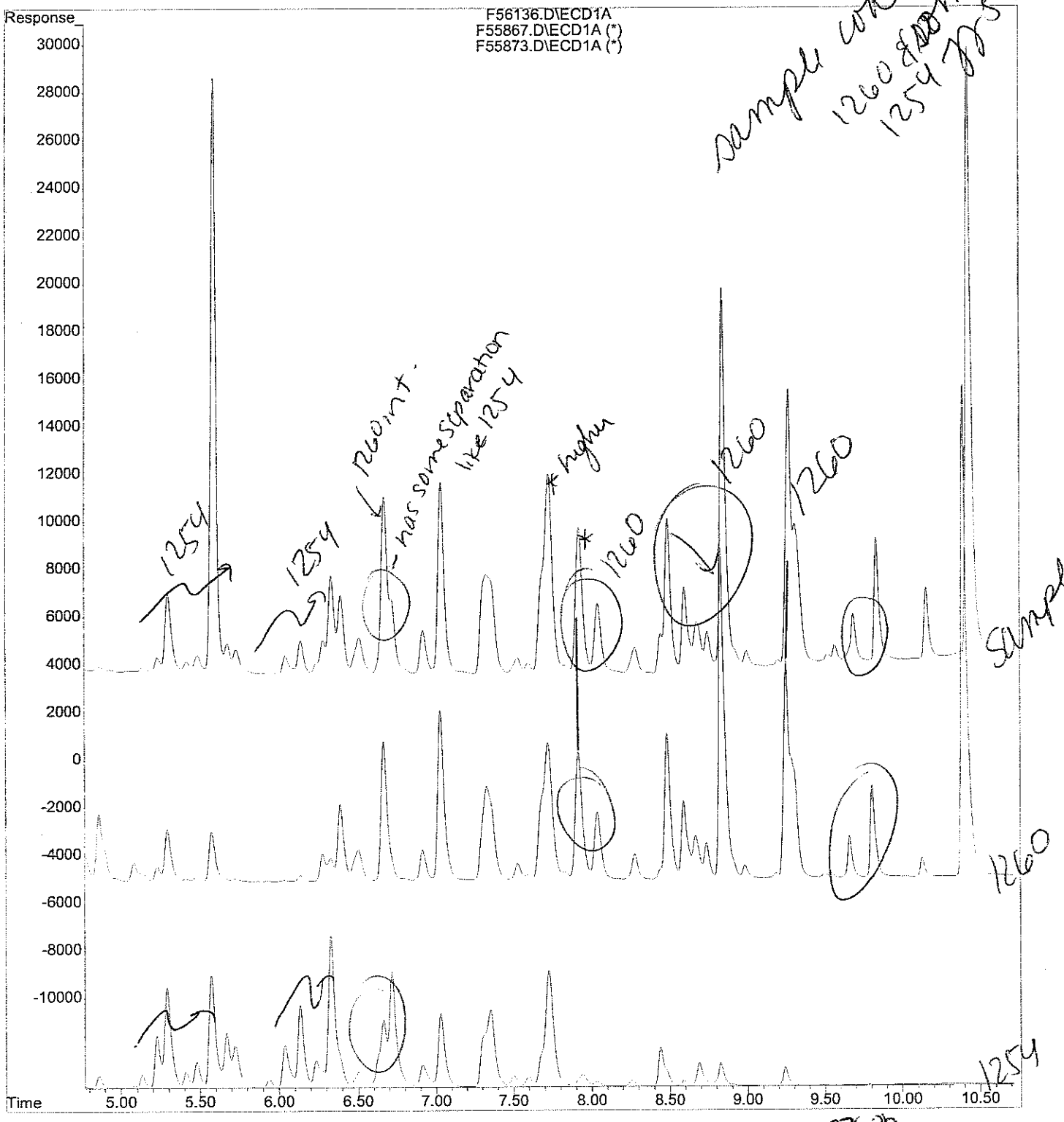
Signal #1 : C:\HPCHEM\2\DATA\052606-F\F56136.D\ECD1A.CH Vial: 53
Signal #2 : C:\HPCHEM\2\DATA\052606-F\F56136.D\ECD2B.CH
Acq On : 26 May 2006 2:51 pm Operator:
Sample : 56617-3, 1:5, A/C Inst : Instrumen
Misc : SOIL Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 26 15:06 2006 Quant Results File: PBQ05166.RES

Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 10:06:32 2006
Response via : Multiple Level Calibration
DataAcq Meth : PSTQUICK.M

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film



File : C:\HPCHEM\2\DATA\052606-F\F56136.D
Operator :
Acquired : 26 May 2006 2:51 pm using AcqMethod PSTQUICK.M
Instrument : Instrumen
Sample Name: 56617-3, 1:5, A/C
Misc Info : SOIL
Vial Number: 53



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May 26, 2006

CLIENT SAMPLE ID

Project Name: GP

Project Number: 218736

Field Sample ID: SB-10 (0-1)

SAMPLE DATA

Lab Sample ID: 56617-4
Matrix: Solid
Percent Solid: 97
Dilution Factor: 10
Collection Date: 05/24/06
Lab Receipt Date: 05/25/06
Extraction Date: 05/25/06
Analysis Date: 05/26/06

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit µg/kg	Results µg/kg
PCB-1016	150	U
PCB-1221	150	U
PCB-1232	150	U
PCB-1242	150	**1290**
PCB-1248	150	U
PCB-1254	150	U
PCB-1260	150	U

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	90	%
Decachlorobiphenyl	90	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082.

Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3545.

COMMENTS: Results are expressed on a dry weight basis.

PCB Report

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PCB CONFIRMATION RESULTS

SDG: 56617

Project No: Site ID: Location: ME

Instrument ID: F GC Columns: RTX-CLPesticidesI ID: 0.32(mm)

RTX-CLPesticidesII

Sample ID: 56617-4

COMPOUND	RESPONSE #1 ug/kg	RESPONSE #2 ug/kg	DIFFERENCE %	#	% DIFFERENCE Acceptance Limit
PCB 1242	1290	1420	10		40

Column to be used to flag % Difference with an asterisk

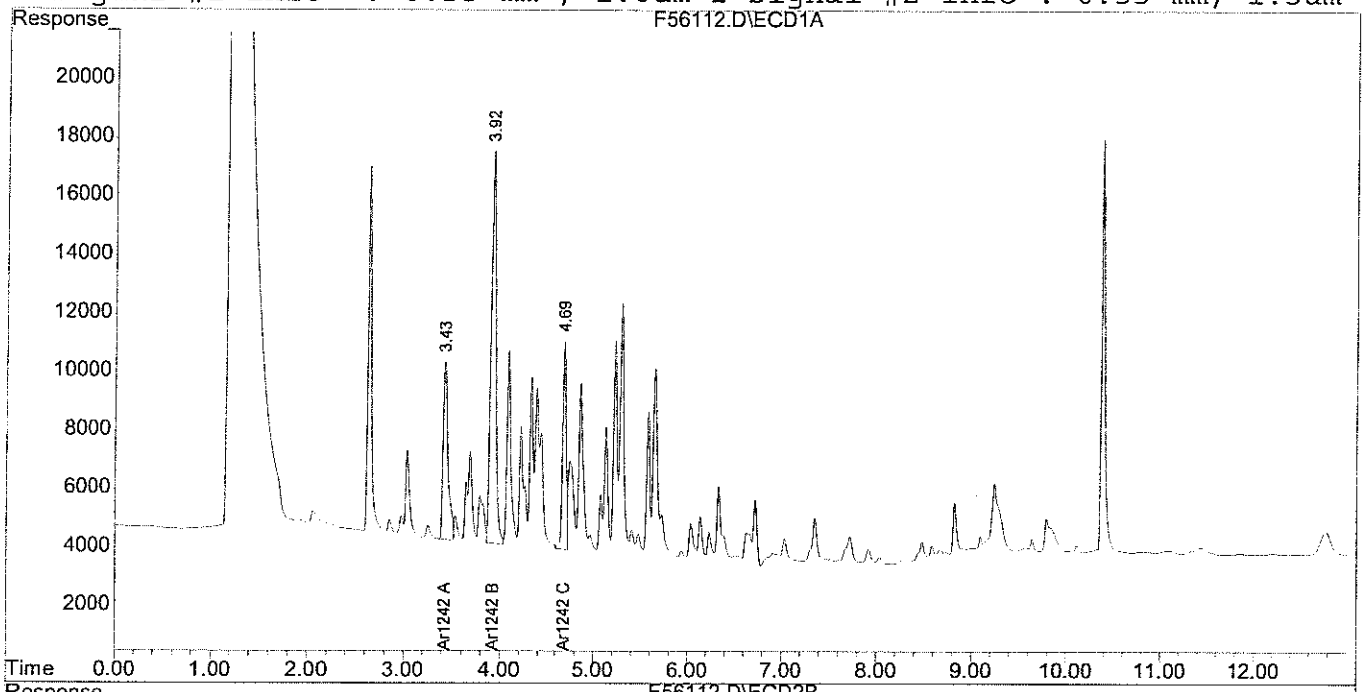
* Values outside of QC limits

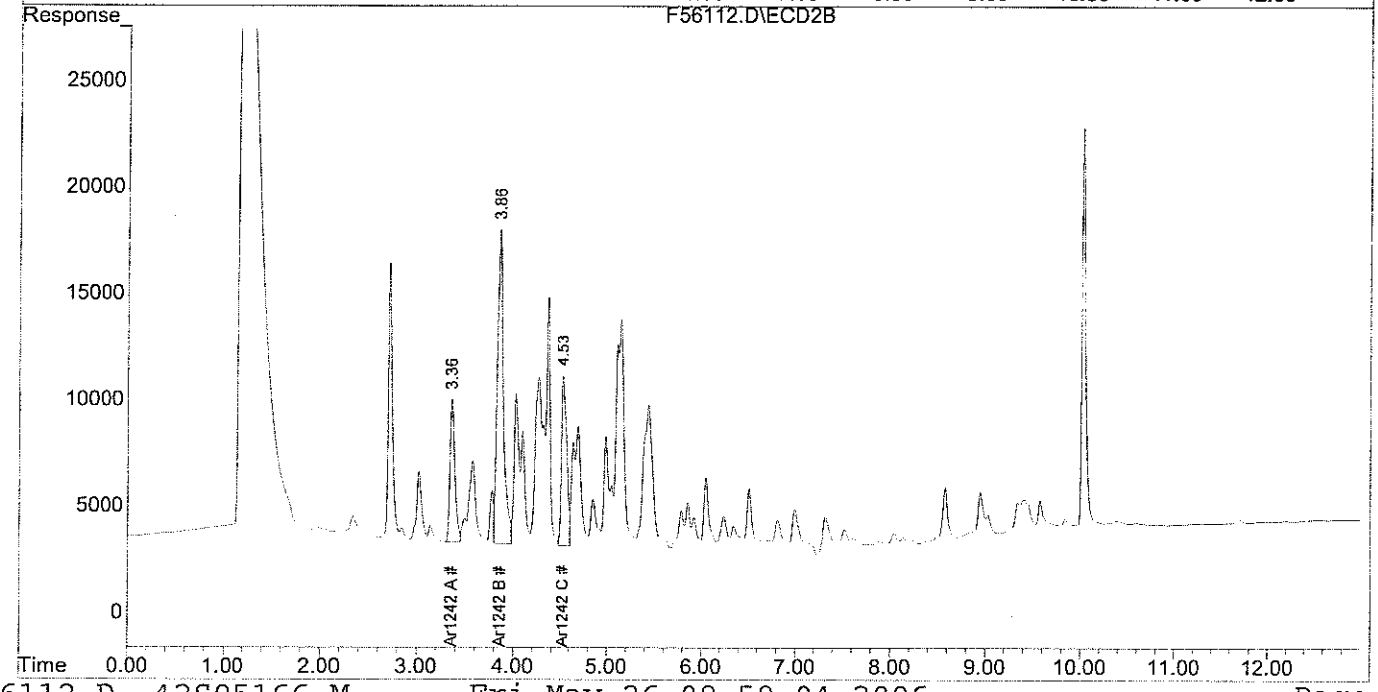
Quantitation Report (QT Reviewed)

Signal #1 : C:\HPCHEM\2\DATA\052506-F\F56112.D\ECD1A.CH Vial: 34
Signal #2 : C:\HPCHEM\2\DATA\052506-F\F56112.D\ECD2B.CH
Acq On : 26 May 2006 12:26 am Operator:
Sample : 56617-4, 1:10, A/C Inst : Instrumen
Misc : SOIL Multiplr: 1.00
IntFile Signal #1: EVENTS.E IntFile Signal #2: EVENTS2.E
Quant Time: May 26 9:58 2006 Quant Results File: 42S05166.RES

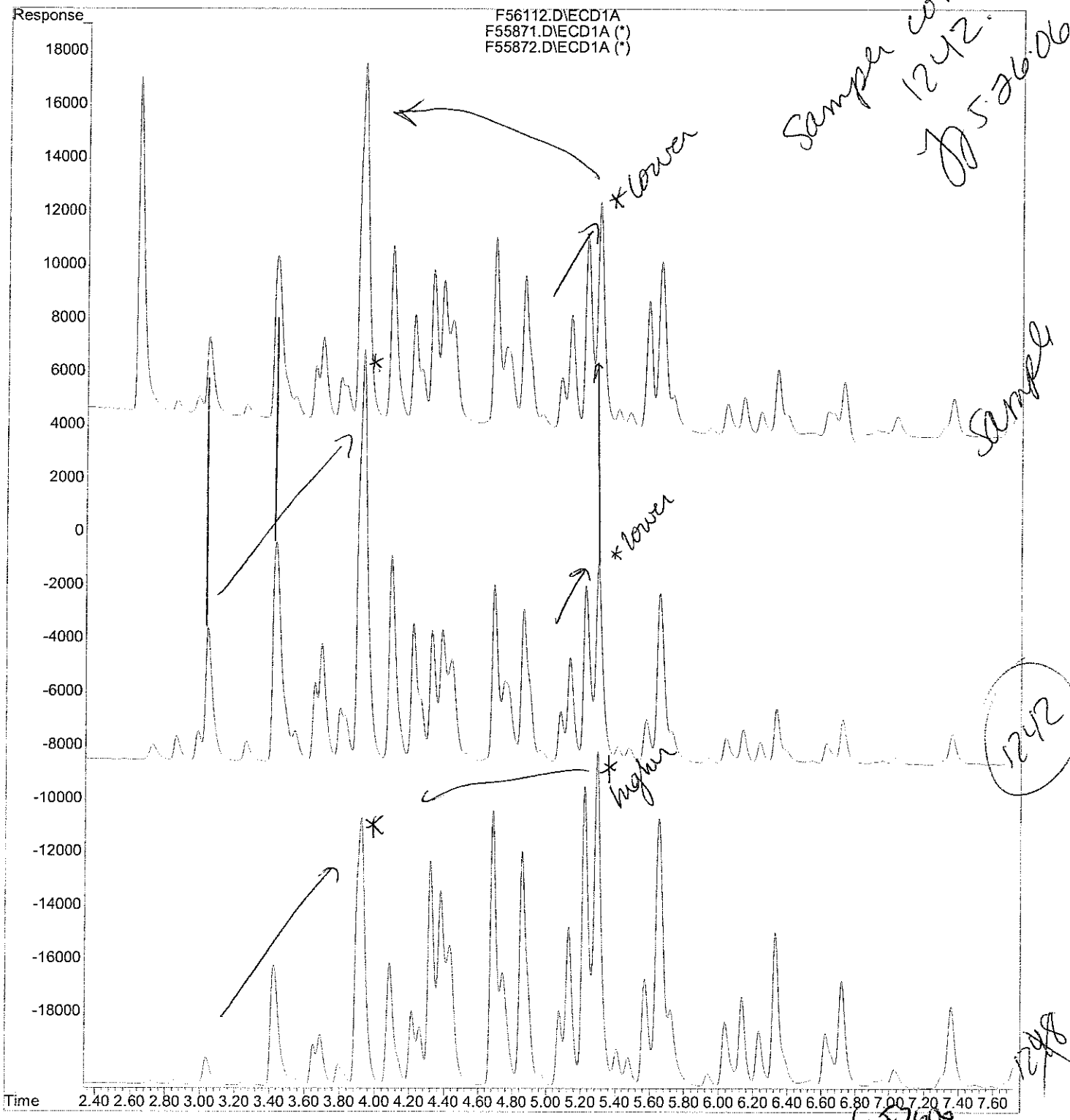
Quant Method : C:\HPCHEM\2\METHODS\42S05166.M (Chemstation Integrator)
Title : Aroclor 1242
Last Update : Fri May 26 08:34:45 2006
Response via : Single Level Calibration
DataAcq Meth : PSTQUICK.M

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film





File : C:\HPCHEM\2\DATA\052506-F\F56112.D
Operator :
Acquired : 26 May 2006 12:26 am using AcqMethod PSTQUICK.M
Instrument : Instrumen
Sample Name: 56617-4, 1:10, A/C
Misc Info : SOIL
Vial Number: 34



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May 26, 2006

CLIENT SAMPLE ID

Project Name: GP

Project Number: 218736

Field Sample ID: SB-11 (0-4)

SAMPLE DATA

Lab Sample ID: 56617-5
Matrix: Solid
Percent Solid: 93
Dilution Factor: 1.1
Collection Date: 05/24/06
Lab Receipt Date: 05/25/06
Extraction Date: 05/25/06
Analysis Date: 05/26/06

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit µg/kg	Results µg/kg
PCB-1016	17	**78***
PCB-1221	17	U
PCB-1232	17	U
PCB-1242	17	U
PCB-1248	17	U
PCB-1254	17	U
PCB-1260	17	**72**

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	99	%
Decachlorobiphenyl	79	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082.

COMMENTS: Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3545. Results are expressed on a dry weight basis. *See narrative.

PCB Report

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PCB CONFIRMATION RESULTS

SDG: 56617

Project No: Site ID: Location: ME

Instrument ID: F GC Columns: RTX-CLPesticidesI ID: 0.32(mm)
RTX-CLPesticidesII

Sample ID: 56617-5

COMPOUND	RESPONSE #1 ug/kg	RESPONSE #2 ug/kg	DIFFERENCE %	#	% DIFFERENCE Acceptance Limit
PCB 1016	78	56	33		40
PCB 1260	71	89	22		40

Column to be used to flag % Difference with an asterisk

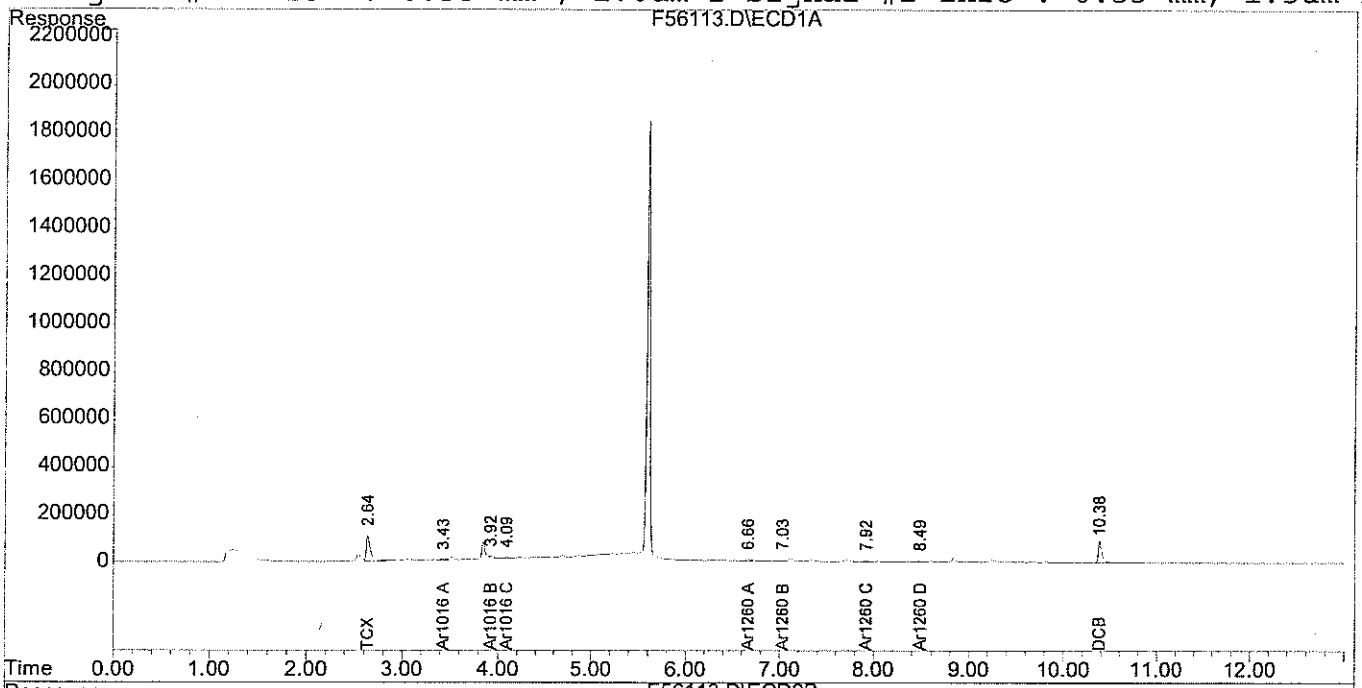
* Values outside of QC limits

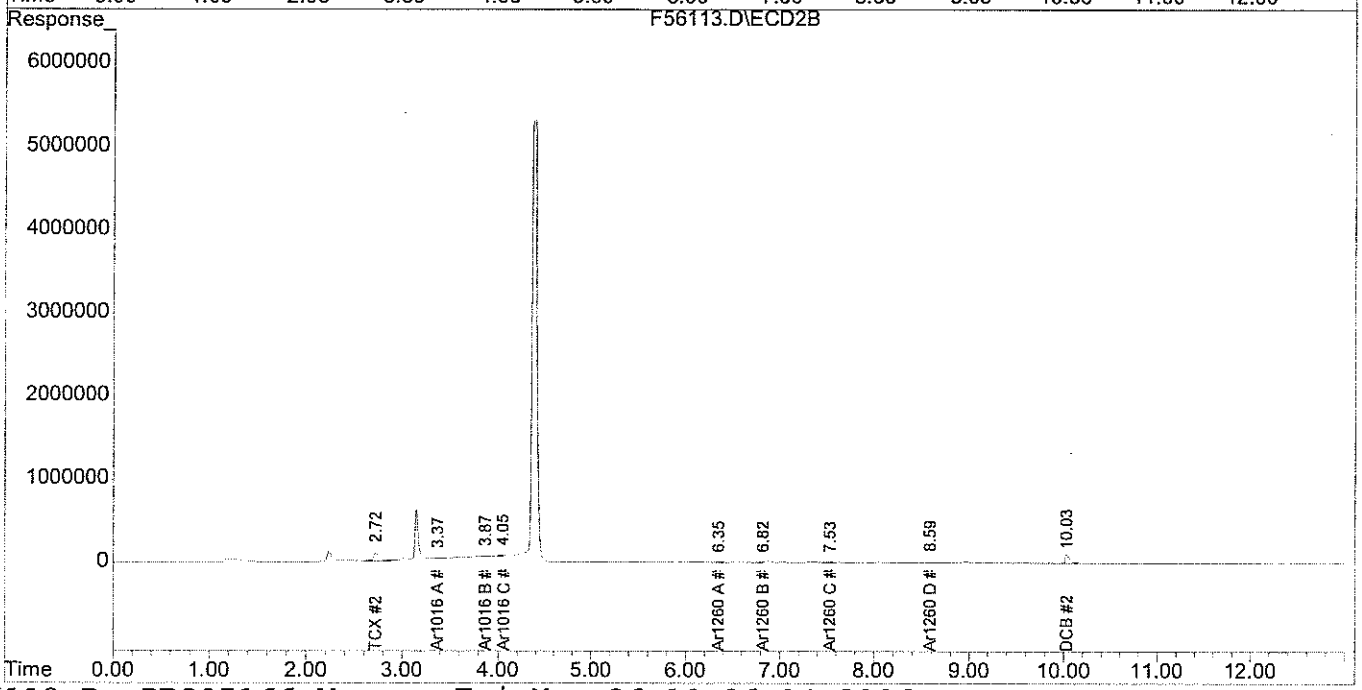
Quantitation Report (QT Reviewed)

Signal #1 : C:\HPCHEM\2\DATA\052506-F\F56113.D\ECD1A.CH Vial: 35
Signal #2 : C:\HPCHEM\2\DATA\052506-F\F56113.D\ECD2B.CH
Acq On : 26 May 2006 12:41 am Operator:
Sample : 56617-5, A/C Inst : Instrumen
Misc : SOIL Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 26 10:25 2006 Quant Results File: PBQ05166.RES

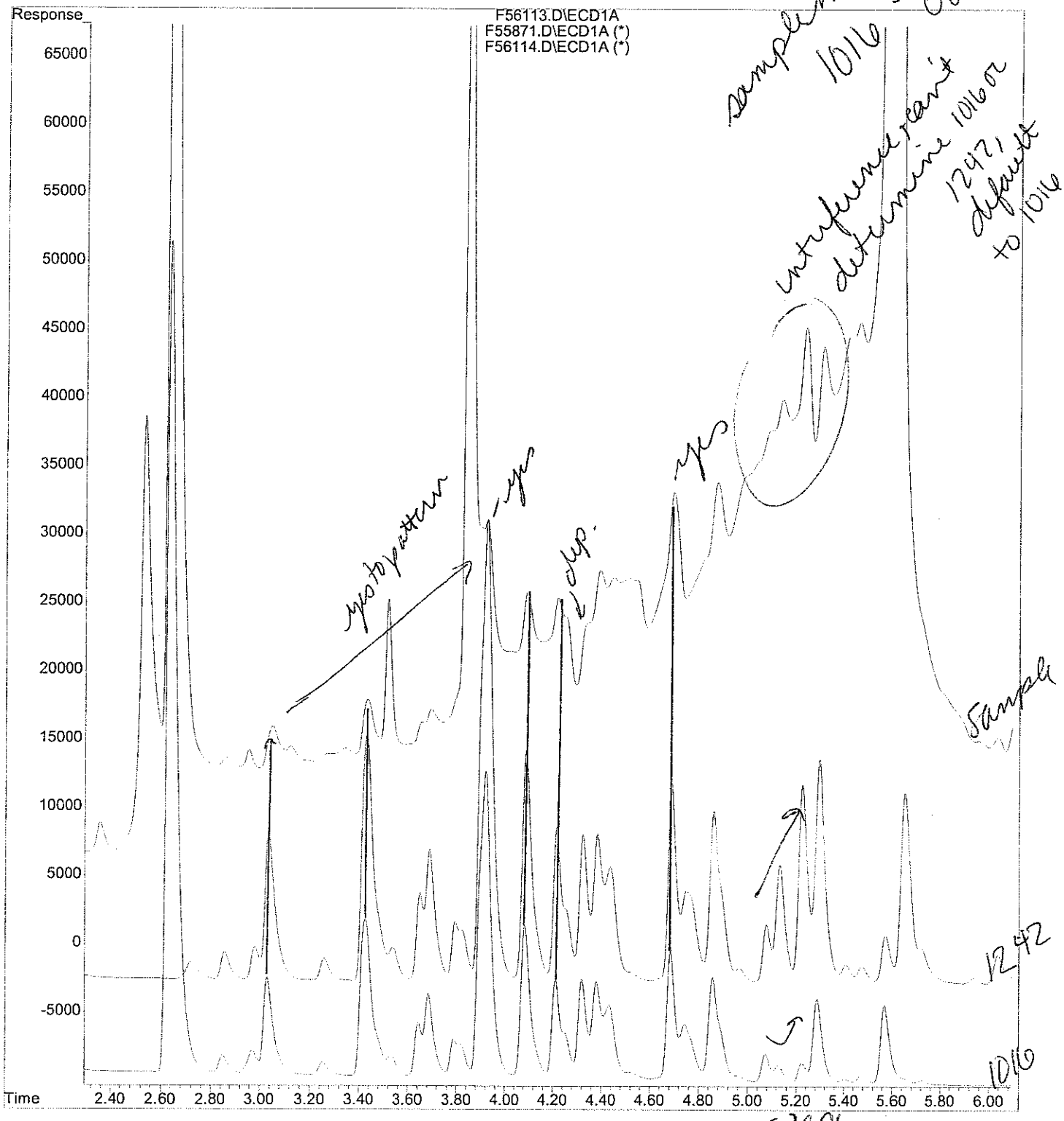
Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 10:06:32 2006
Response via : Multiple Level Calibration
DataAcq Meth : PSTQUICK.M

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film

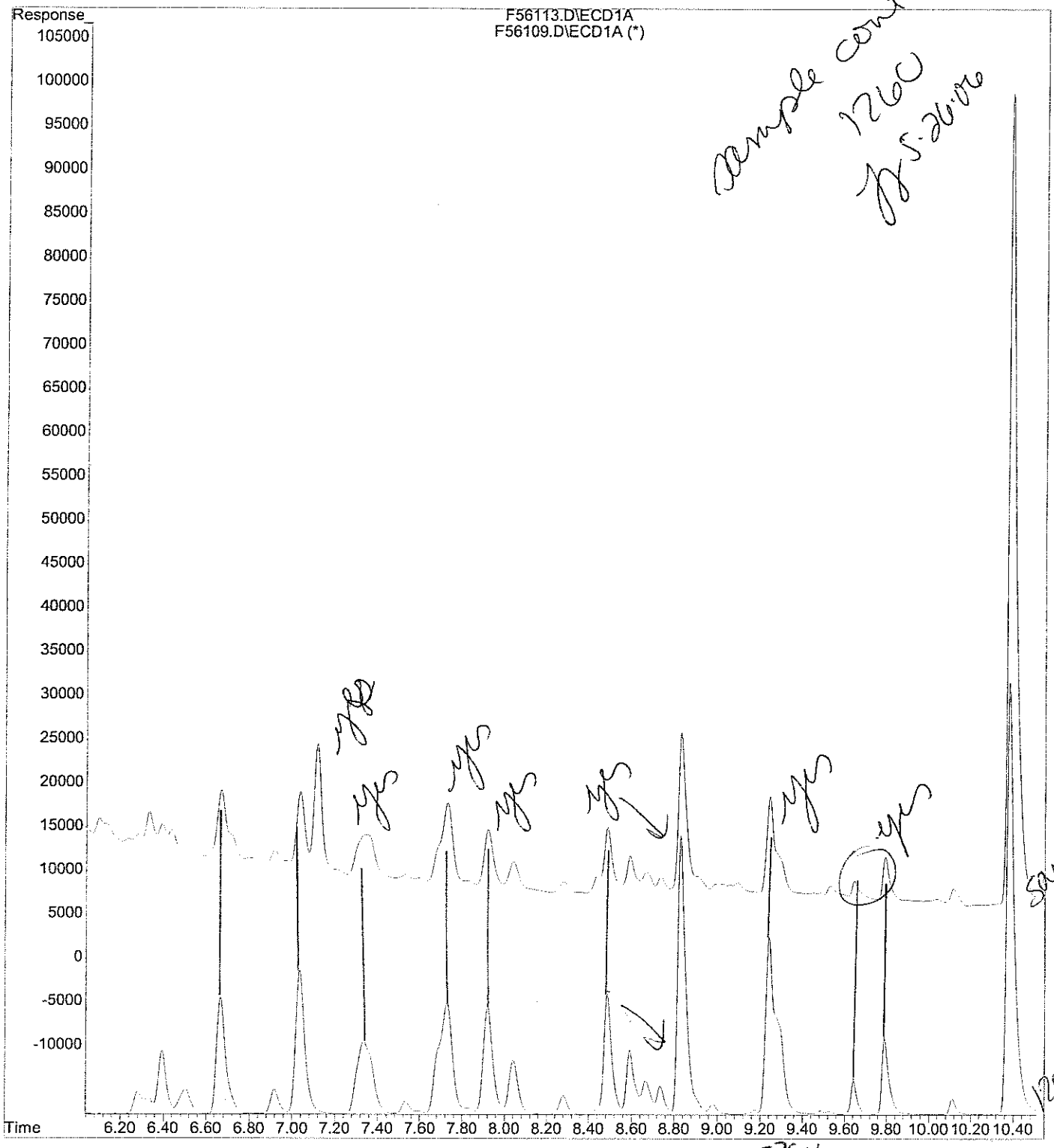




File : C:\HPCHEM\2\DATA\052506-F\F56113.D
Operator :
Acquired : 26 May 2006 12:41 am using AcqMethod PSTQUICK.M
Instrument : Instrumen
Sample Name: 56617-5, A/C
Misc Info : SOIL
Vial Number: 35



File : C:\HPCHEM\2\DATA\052506-F\F56113.D
Operator :
Acquired : 26 May 2006 12:41 am using AcqMethod PSTQUICK.M
Instrument : Instrumen
Sample Name: 56617-5, A/C
Misc Info : SOIL
Vial Number: 35



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May 26, 2006

CLIENT SAMPLE ID

Project Name: GP

Project Number: 218736

Field Sample ID: SB-12 (0-4)

SAMPLE DATA

Lab Sample ID: 56617-6
Matrix: Solid
Percent Solid: 92
Dilution Factor: 1.1
Collection Date: 05/24/06
Lab Receipt Date: 05/25/06
Extraction Date: 05/25/06
Analysis Date: 05/25/06

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit µg/kg	Results µg/kg
PCB-1016	17	U
PCB-1221	17	U
PCB-1232	17	U
PCB-1242	17	U
PCB-1248	17	U
PCB-1254	17	U
PCB-1260	17	**51**

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	78	%
Decachlorobiphenyl	83	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082.

Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3545.

COMMENTS: Results are expressed on a dry weight basis.

PCB Report

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PCB CONFIRMATION RESULTS

SDG: 56617

Project No: Site ID: Location: ME

Instrument ID: F GC Columns: RTX-CLPesticidesI ID: 0.32(mm)

RTX-CLPesticidesII

Sample ID: 56617-6

COMPOUND	RESPONSE #1 ug/kg	RESPONSE #2 ug/kg	DIFFERENCE %	#	% DIFFERENCE Acceptance Limit
PCB 1260	51	49	4		40

Column to be used to flag % Difference with an asterisk

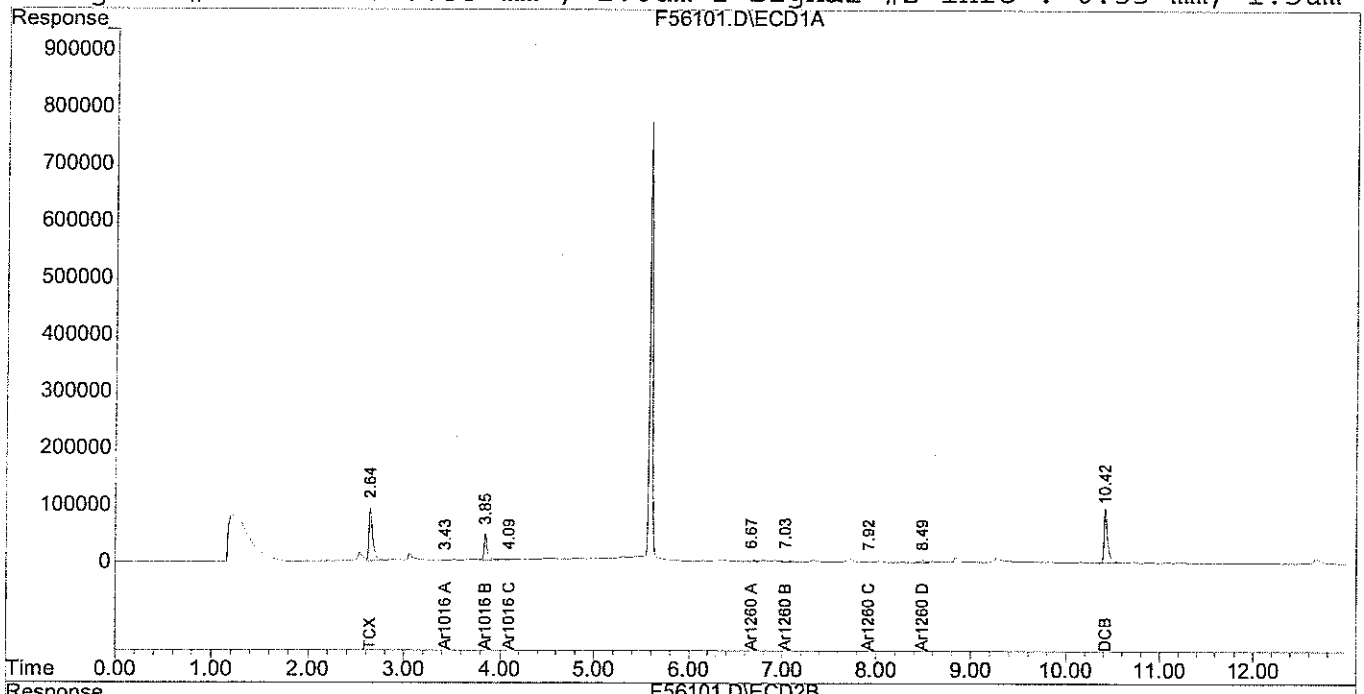
* Values outside of QC limits

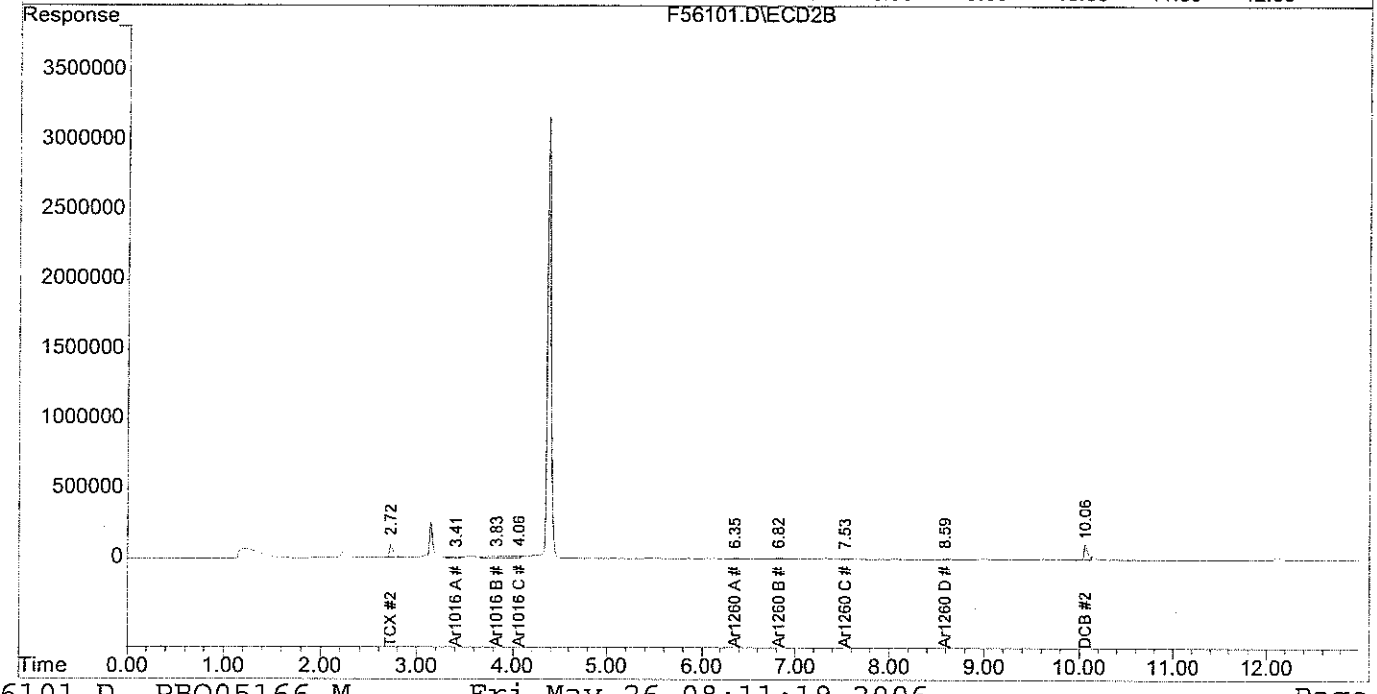
Quantitation Report (QT Reviewed)

Signal #1 : C:\HPCHEM\2\DATA\052506-F\F56101.D\ECD1A.CH Vial: 24
Signal #2 : C:\HPCHEM\2\DATA\052506-F\F56101.D\ECD2B.CH
Acq On : 25 May 2006 9:16 pm Operator:
Sample : 56617-6, A/C Inst : Instrumen
Misc : SOIL Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 26 8:11 2006 Quant Results File: PBQ05166.RES

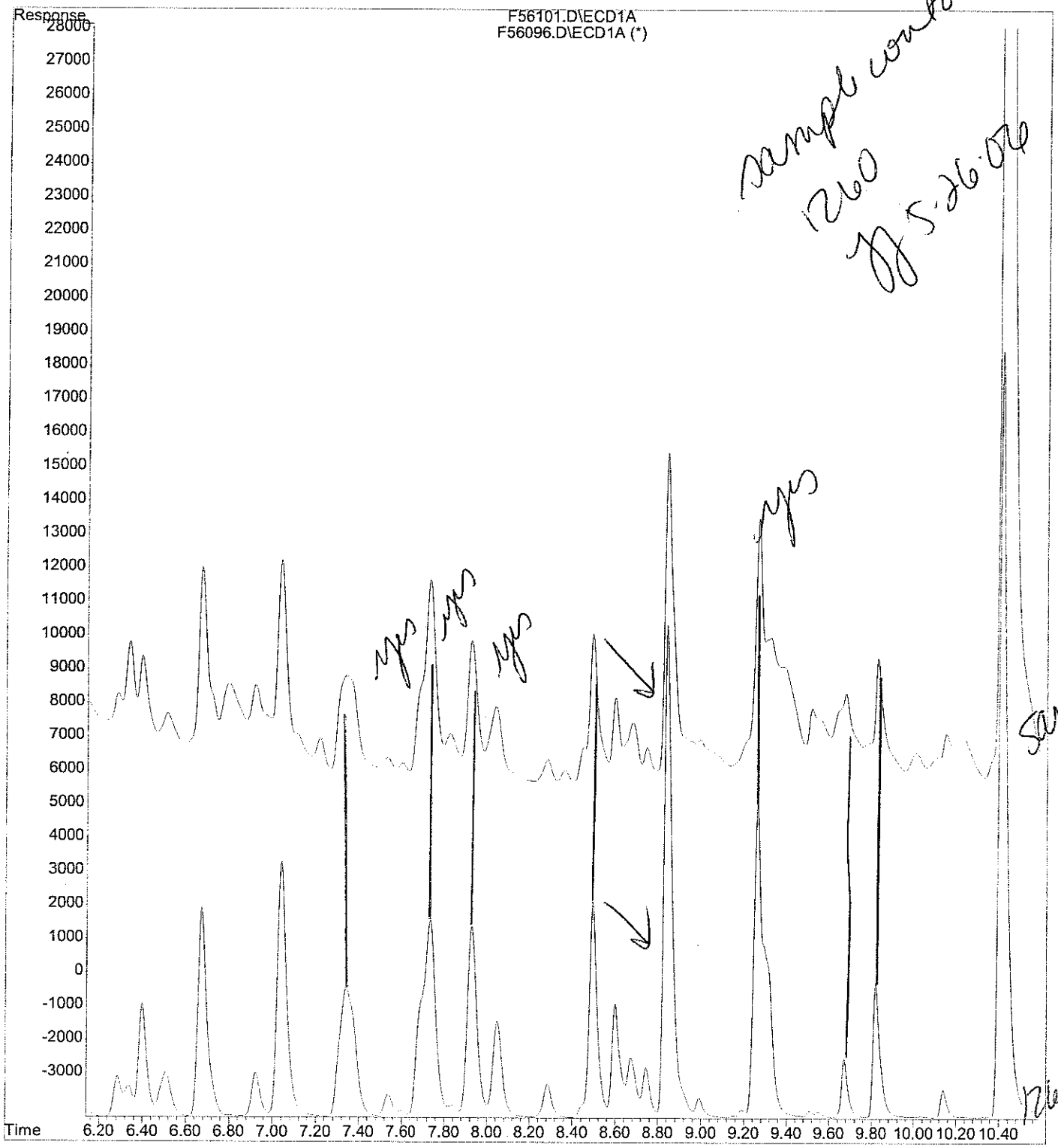
Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 10:06:32 2006
Response via : Multiple Level Calibration
DataAcq Meth : PSTQUICK.M

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film





File : C:\HPCHEM\2\DATA\052506-F\F56101.D
Operator :
Acquired : 25 May 2006 9:16 pm using AcqMethod PSTQUICK.M
Instrument : Instrumen
Sample Name: 56617-6, A/C
Misc Info : SOIL
Vial Number: 24



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May 26, 2006

CLIENT SAMPLE ID

Project Name: GP

Project Number: 218736

Field Sample ID: SB-15 (0-4)

SAMPLE DATA

Lab Sample ID: 56617-7
Matrix: Solid
Percent Solid: 94
Dilution Factor: 1.0
Collection Date: 05/24/06
Lab Receipt Date: 05/25/06
Extraction Date: 05/25/06
Analysis Date: 05/25/06

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit µg/kg	Results µg/kg
PCB-1016	15	U
PCB-1221	15	U
PCB-1232	15	U
PCB-1242	15	U
PCB-1248	15	U
PCB-1254	15	U
PCB-1260	15	U

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	66	%
Decachlorobiphenyl	85	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082.

Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3545.

COMMENTS: Results are expressed on a dry weight basis.

PCB Report

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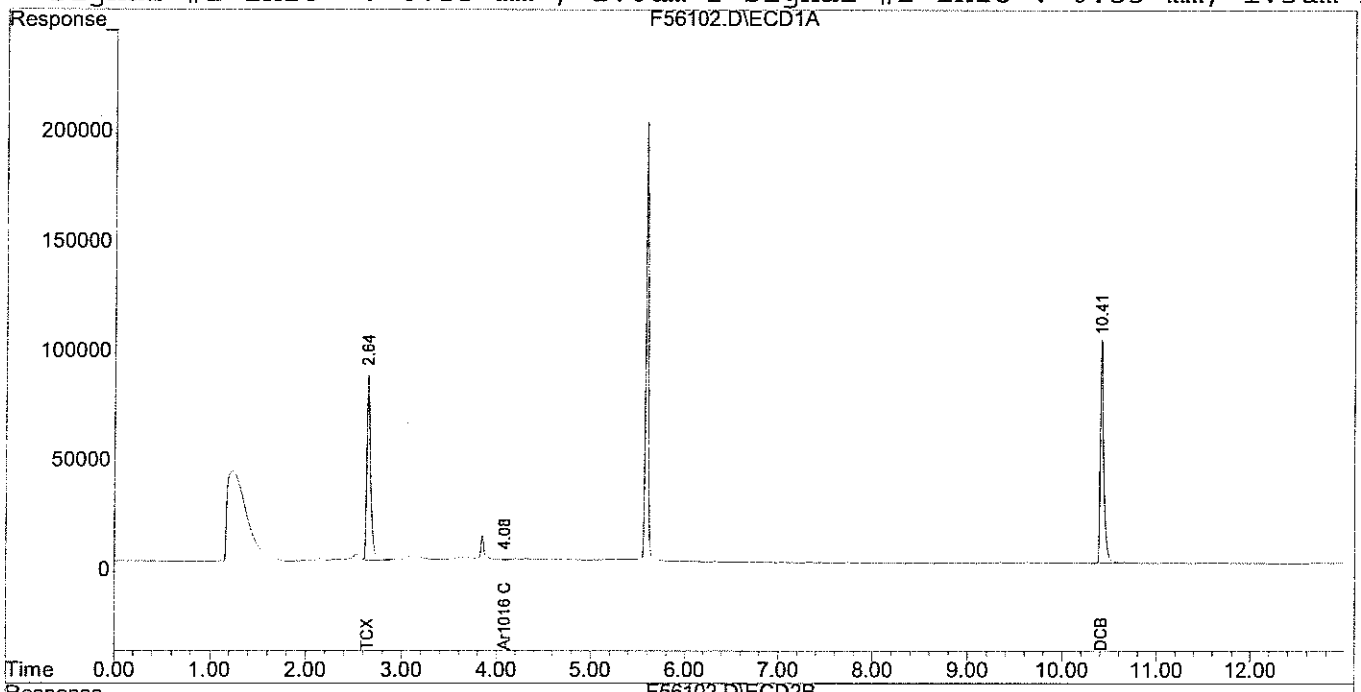
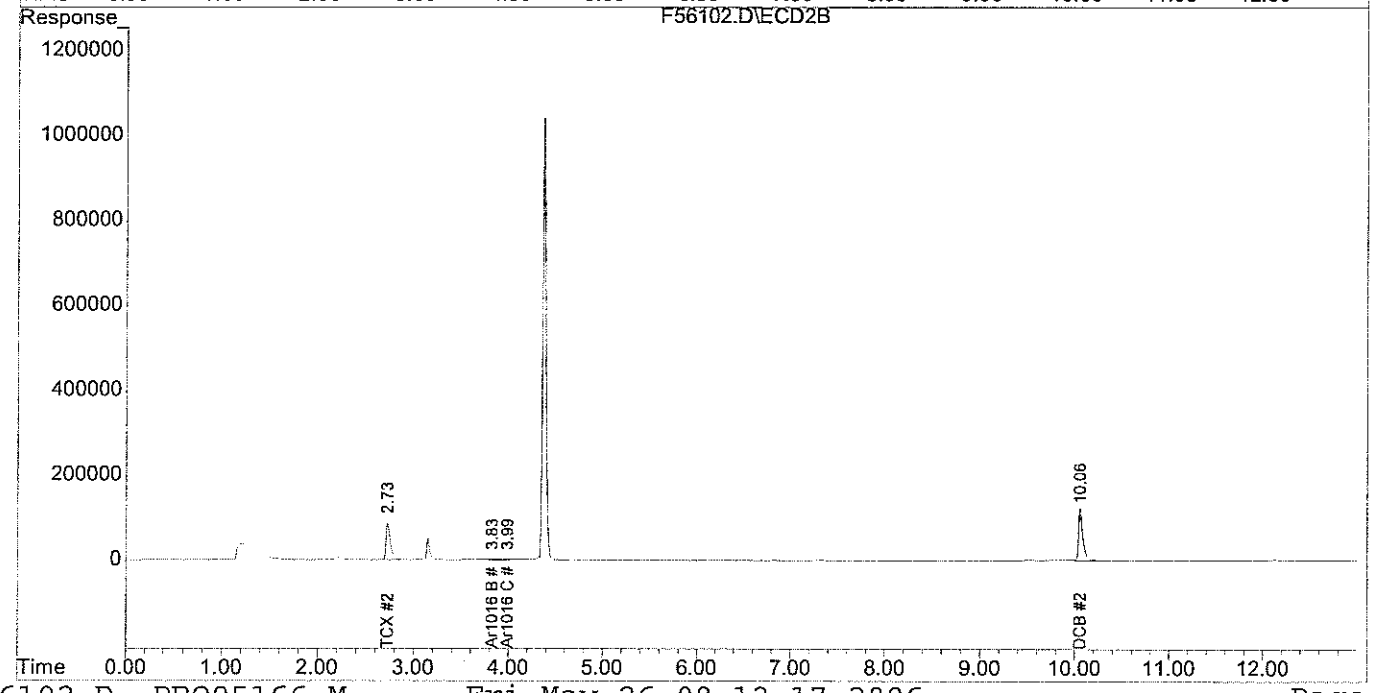
Quantitation Report (QT Reviewed)

Signal #1 : C:\HPCHEM\2\DATA\052506-F\F56102.D\ECD1A.CH Vial: 25
Signal #2 : C:\HPCHEM\2\DATA\052506-F\F56102.D\ECD2B.CH
Acq On : 25 May 2006 9:45 pm Operator:
Sample : 56617-7, A/C Inst : Instrumen
Misc : SOIL Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 26 8:12 2006 Quant Results File: PBQ05166.RES

Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 10:06:32 2006
Response via : Multiple Level Calibration
DataAcq Meth : PSTQUICK.M

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film





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Mr. Karl Kasper
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Portland ME 04102

May 26, 2006

CLIENT SAMPLE ID

Project Name: GP

Project Number: 218736

Field Sample ID: SB-18 (0-4)

SAMPLE DATA

Lab Sample ID: 56617-8
Matrix: Solid
Percent Solid: 88
Dilution Factor: 1.1
Collection Date: 05/24/06
Lab Receipt Date: 05/25/06
Extraction Date: 05/25/06
Analysis Date: 05/25/06

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit µg/kg	Results µg/kg
PCB-1016	17	U
PCB-1221	17	U
PCB-1232	17	U
PCB-1242	17	U
PCB-1248	17	U
PCB-1254	17	U
PCB-1260	17	U

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	69	%
Decachlorobiphenyl	79	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082.

Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3545.

COMMENTS: Results are expressed on a dry weight basis.

PCB Report

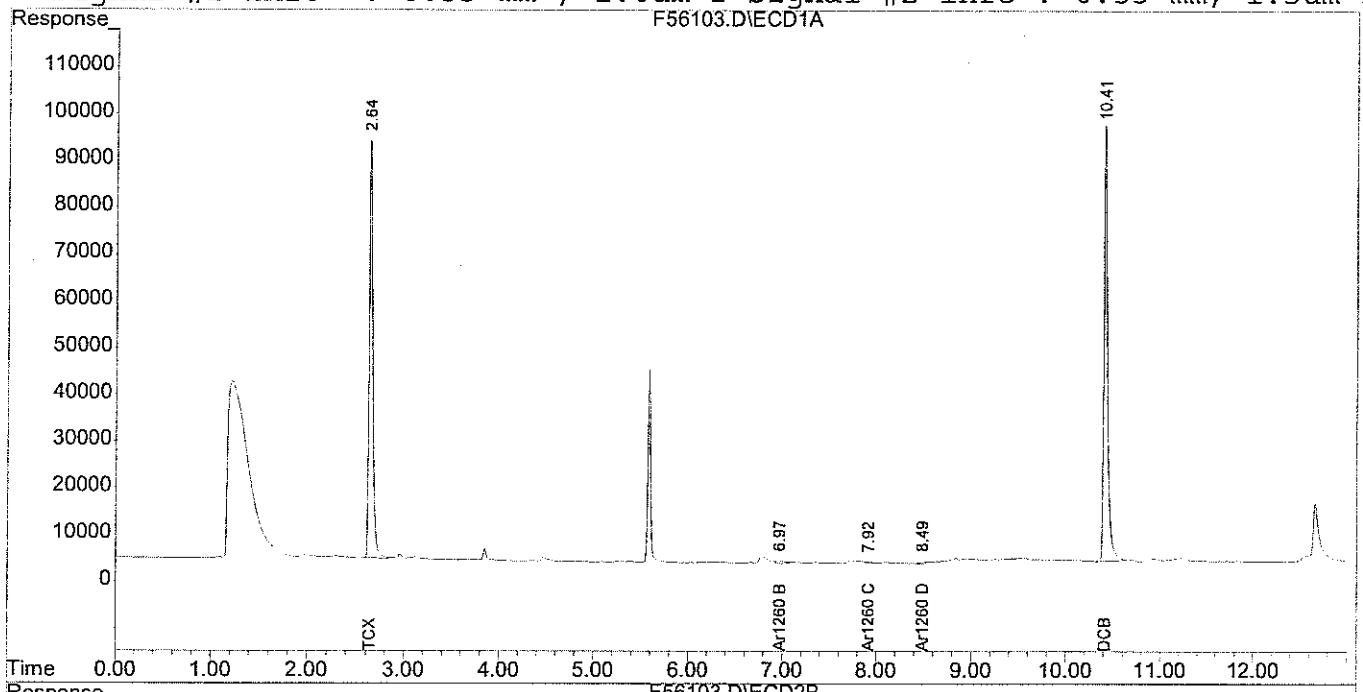
Authorized signature Milena Tull

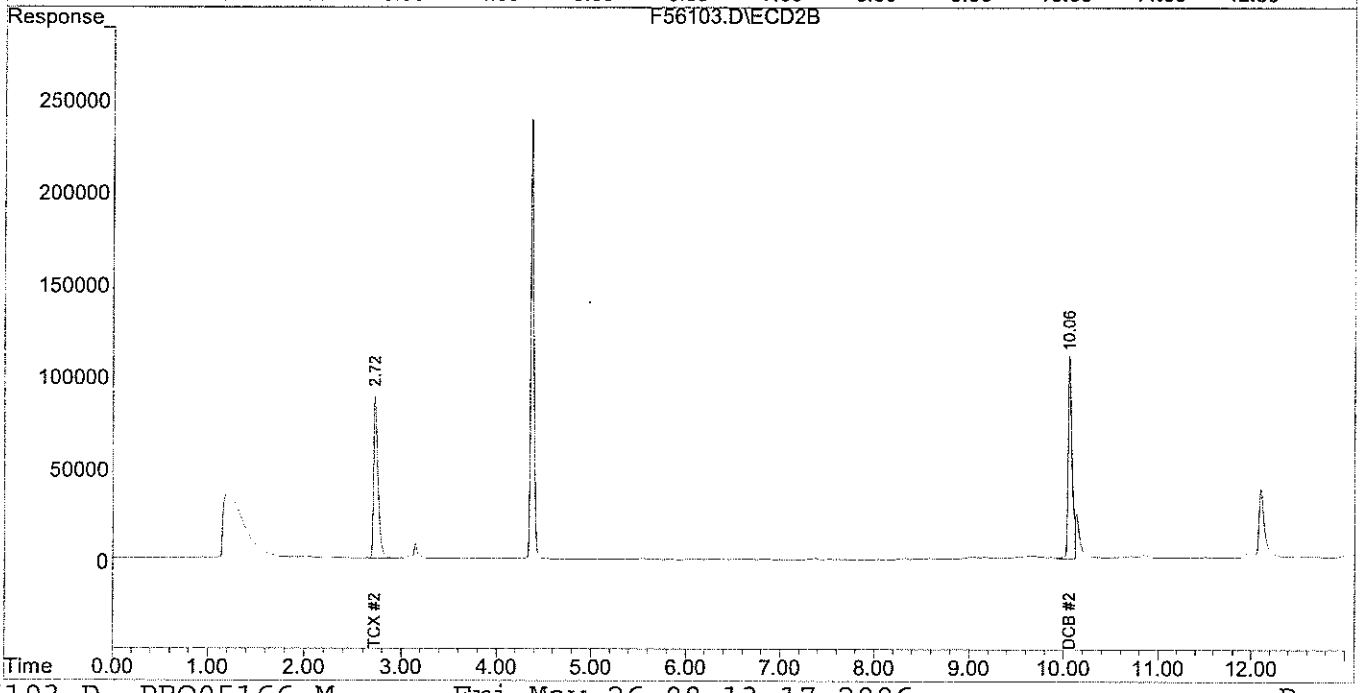
Quantitation Report (QT Reviewed)

Signal #1 : C:\HPCHEM\2\DATA\052506-F\F56103.D\ECD1A.CH Vial: 26
Signal #2 : C:\HPCHEM\2\DATA\052506-F\F56103.D\ECD2B.CH
Acq On : 25 May 2006 10:00 pm Operator:
Sample : 56617-8, A/C Inst : Instrumen
Misc : SOIL Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 26 8:13 2006 Quant Results File: PBQ05166.RES

Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 10:06:32 2006
Response via : Multiple Level Calibration
DataAcq Meth : PSTQUICK.M

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film





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May 26, 2006

CLIENT SAMPLE ID

Project Name: GP

Project Number: 218736

Field Sample ID: SB-25 (0-4)

SAMPLE DATA

Lab Sample ID: 56617-9
Matrix: Solid
Percent Solid: 96
Dilution Factor: 1.0
Collection Date: 05/24/06
Lab Receipt Date: 05/25/06
Extraction Date: 05/25/06
Analysis Date: 05/25/06

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit µg/kg	Results µg/kg
PCB-1016	15	U
PCB-1221	15	U
PCB-1232	15	U
PCB-1242	15	U
PCB-1248	15	U
PCB-1254	15	U
PCB-1260	15	77

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	84	%
Decachlorobiphenyl	86	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082.

Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3545.

COMMENTS: Results are expressed on a dry weight basis.

PCB Report

Authorized signature

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PCB CONFIRMATION RESULTS

SDG: 56617

Project No: Site ID: Location: ME

Instrument ID: F GC Columns: RTX-CLPesticidesI ID: 0.32(mm)

RTX-CLPesticidesII

Sample ID: 56617-9

COMPOUND	RESPONSE #1 ug/kg	RESPONSE #2 ug/kg	DIFFERENCE %	#	% DIFFERENCE Acceptance Limit
PCB 1260	77	94	20		40

Column to be used to flag % Difference with an asterisk

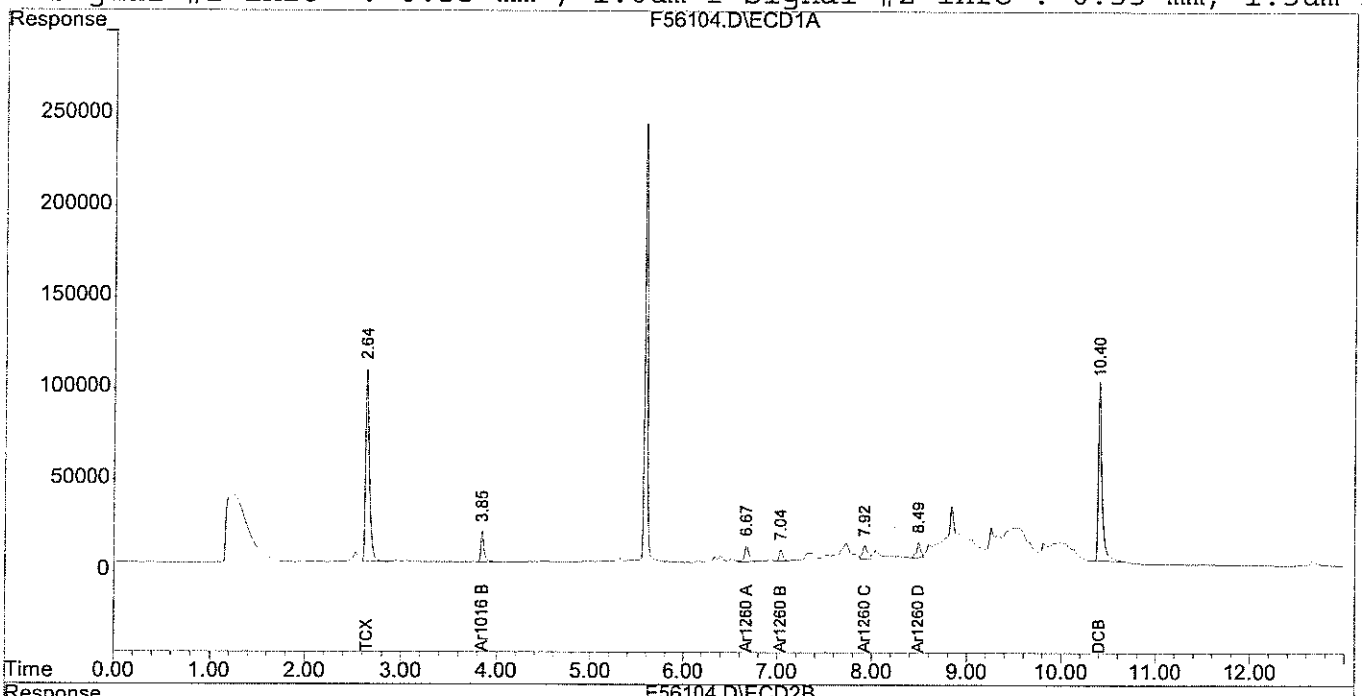
* Values outside of QC limits

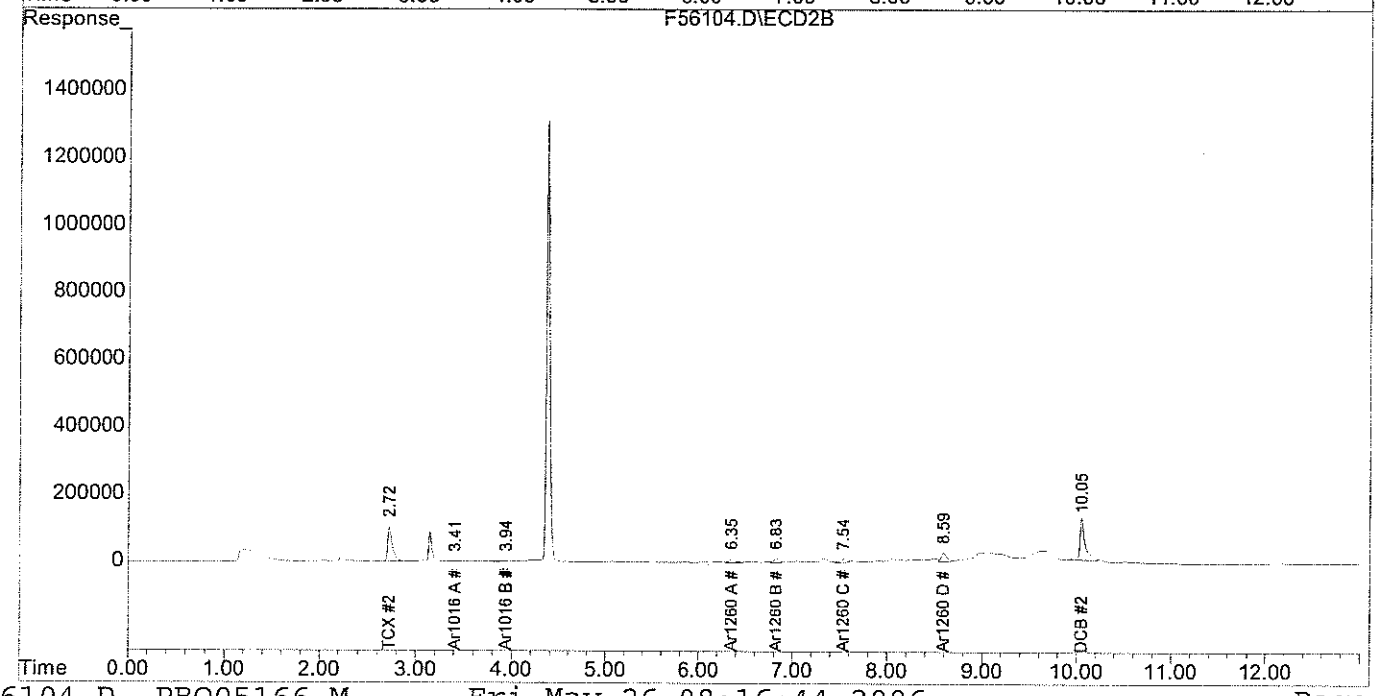
Quantitation Report (QT Reviewed)

Signal #1 : C:\HPCHEM\2\DATA\052506-F\F56104.D\ECD1A.CH Vial: 27
Signal #2 : C:\HPCHEM\2\DATA\052506-F\F56104.D\ECD2B.CH
Acq On : 25 May 2006 10:14 pm Operator:
Sample : 56617-9, A/C Inst : Instrumen
Misc : SOIL Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 26 8:16 2006 Quant Results File: PBQ05166.RES

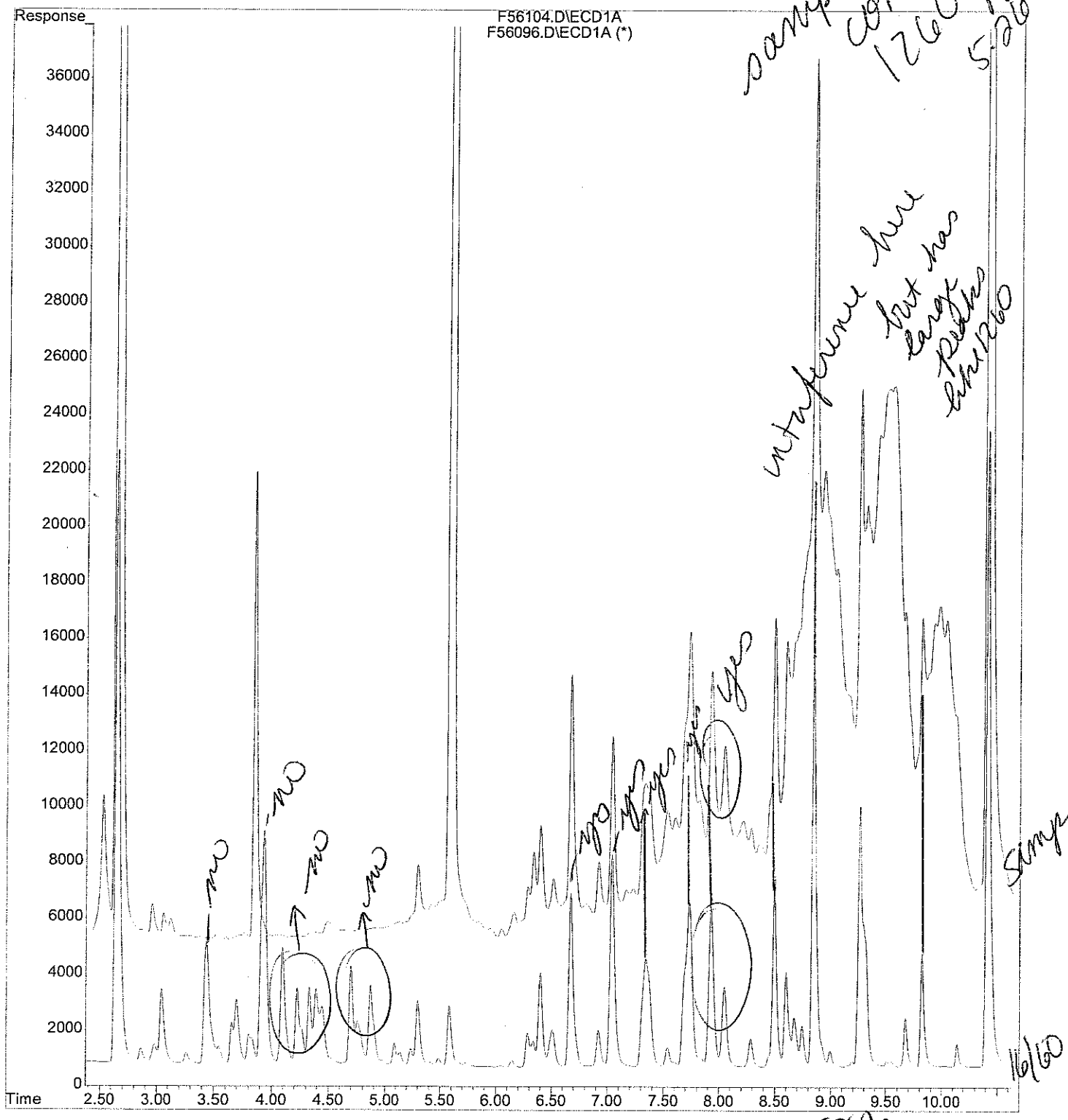
Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 10:06:32 2006
Response via : Multiple Level Calibration
DataAcq Meth : PSTQUICK.M

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film





File : C:\HPCHEM\2\DATA\052506-F\F56104.D
Operator :
Acquired : 25 May 2006 10:14 pm using AcqMethod PSTQUICK.M
Instrument : Instrumen
Sample Name: 56617-9, A/C
Misc Info : SOIL
Vial Number: 27



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May 26, 2006

CLIENT SAMPLE ID

Project Name: GP

Project Number: 218736

Field Sample ID: SB-26 (0-4)

SAMPLE DATA

Lab Sample ID: 56617-10
Matrix: Solid
Percent Solid: 94
Dilution Factor: 1.0
Collection Date: 05/24/06
Lab Receipt Date: 05/25/06
Extraction Date: 05/25/06
Analysis Date: 05/25/06

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit µg/kg	Results µg/kg
PCB-1016	15	U
PCB-1221	15	U
PCB-1232	15	U
PCB-1242	15	U
PCB-1248	15	U
PCB-1254	15	U
PCB-1260	15	455

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	75	%
Decachlorobiphenyl	86	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082.

Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3545.

COMMENTS: Results are expressed on a dry weight basis.

PCB Report

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PCB CONFIRMATION RESULTS

SDG: 56617

Project No: Site ID: Location: ME

Instrument ID: F GC Columns: RTX-CLPesticidesI ID: 0.32(mm)

RTX-CLPesticidesII

Sample ID: 56617-10

COMPOUND	RESPONSE #1 ug/kg	RESPONSE #2 ug/kg	DIFFERENCE %	#	% DIFFERENCE Acceptance Limit
PCB 1260	455	470	3		40

Column to be used to flag % Difference with an asterisk

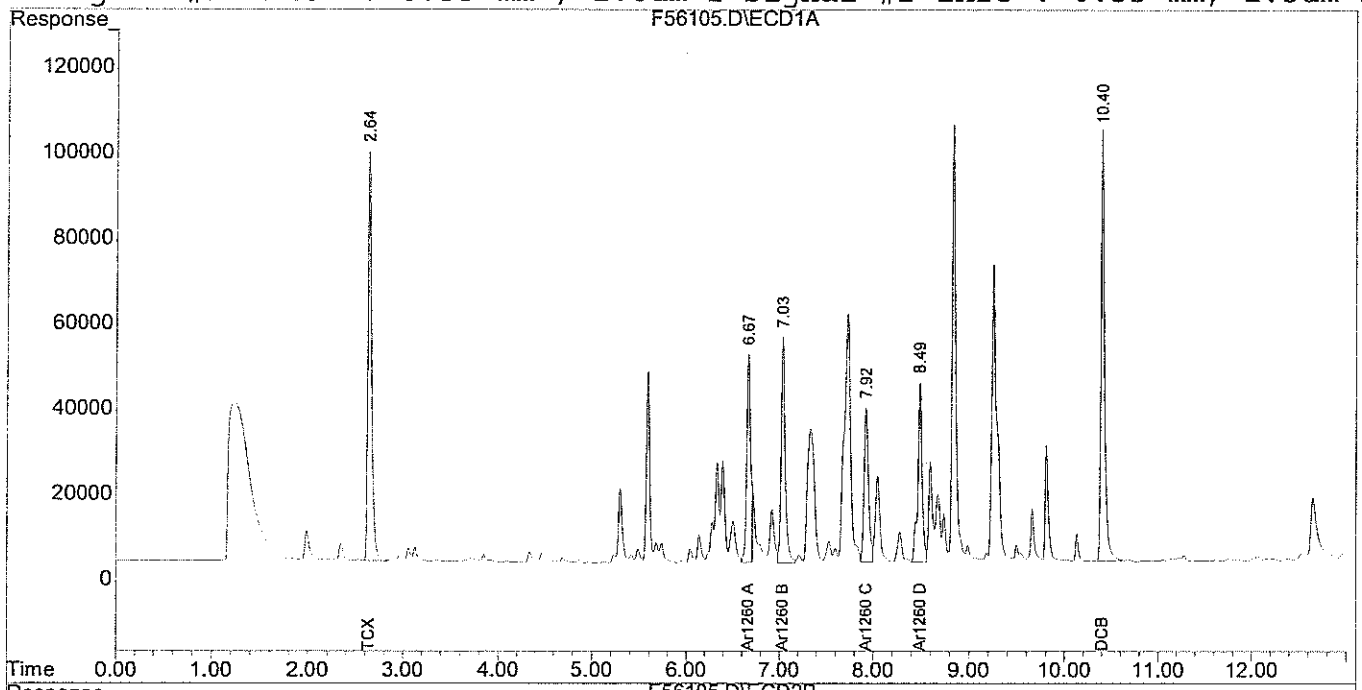
* Values outside of QC limits

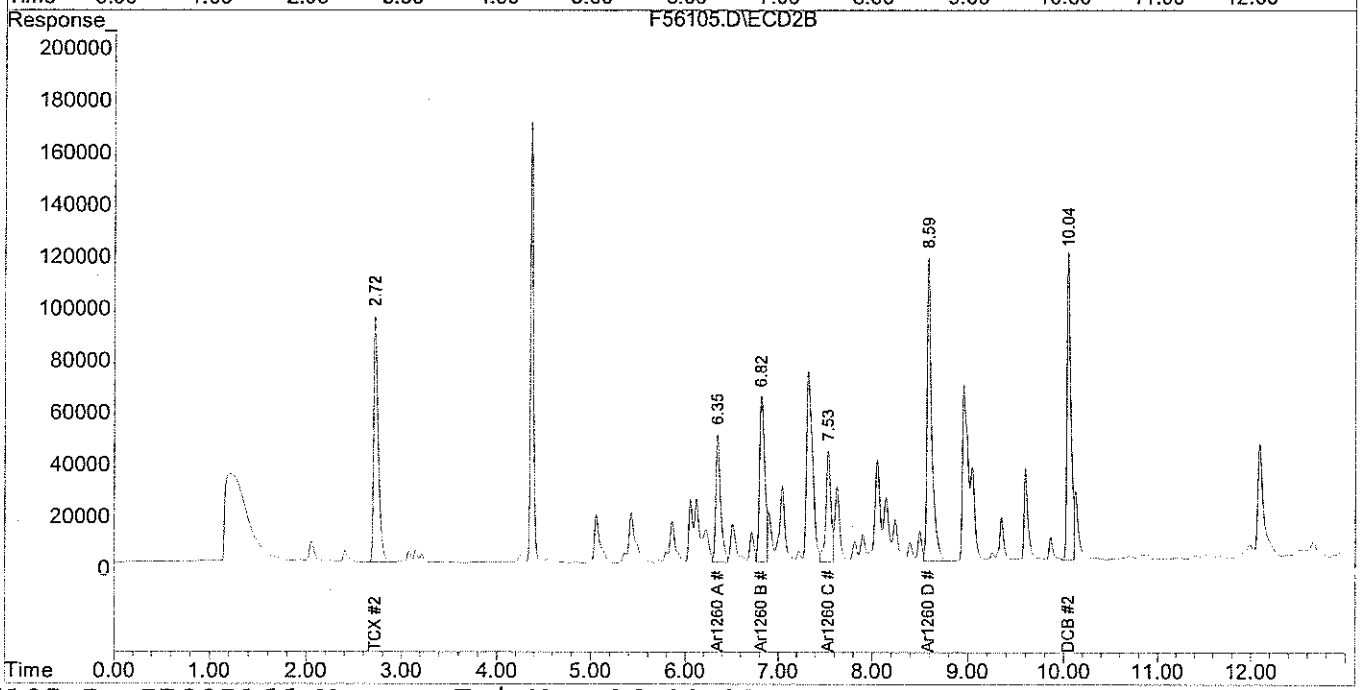
Quantitation Report (QT Reviewed)

Signal #1 : C:\HPCHEM\2\DATA\052506-F\F56105.D\ECD1A.CH Vial: 28
Signal #2 : C:\HPCHEM\2\DATA\052506-F\F56105.D\ECD2B.CH
Acq On : 25 May 2006 10:29 pm Operator:
Sample : 56617-10, A/C Inst : Instrumen
Misc : SOIL Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 26 8:22 2006 Quant Results File: PBQ05166.RES

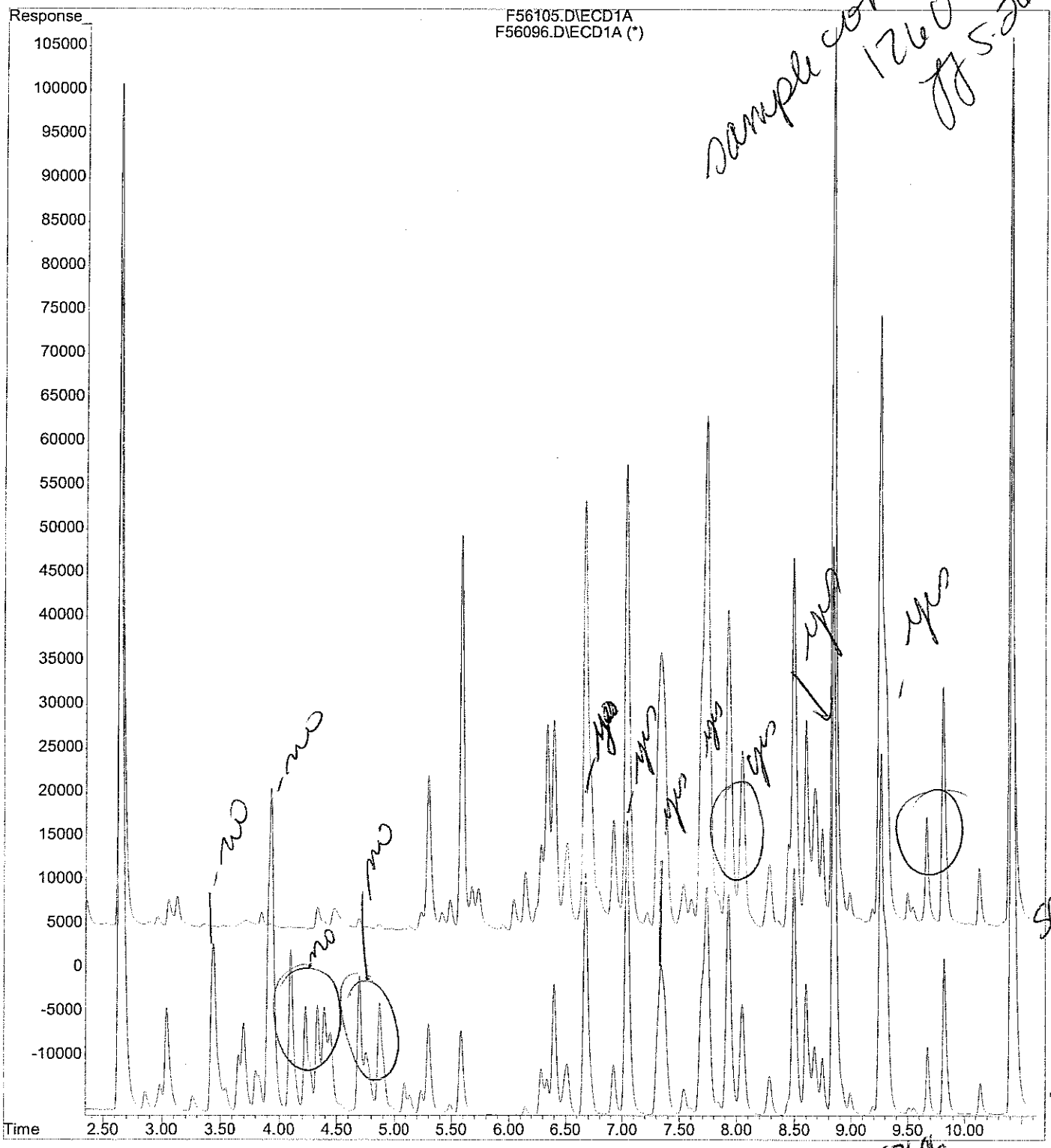
Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 10:06:32 2006
Response via : Multiple Level Calibration
DataAcq Meth : PSTQUICK.M

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film





File : C:\HPCHEM\2\DATA\052506-F\F56105.D
Operator :
Acquired : 25 May 2006 10:29 pm using AcqMethod PSTQUICK.M
Instrument : Instrumen
Sample Name: 56617-10, A/C
Misc Info : SOIL
Vial Number: 28



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Portland ME 04102

May 26, 2006

CLIENT SAMPLE ID

Project Name: GP

Project Number: 218736

Field Sample ID: SB-28 (0-4)

SAMPLE DATA

Lab Sample ID: 56617-11
Matrix: Solid
Percent Solid: 94
Dilution Factor: 1.0
Collection Date: 05/24/06
Lab Receipt Date: 05/25/06
Extraction Date: 05/25/06
Analysis Date: 05/25/06

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit µg/kg	Results µg/kg
PCB-1016	15	U
PCB-1221	15	U
PCB-1232	15	U
PCB-1242	15	U
PCB-1248	15	U
PCB-1254	15	U
PCB-1260	15	**42**

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	79	%
Decachlorobiphenyl	75	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082.

Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3545.

COMMENTS: Results are expressed on a dry weight basis.

PCB Report

Authorized signature Melissa Hall

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PCB CONFIRMATION RESULTS

SDG: 56617

Project No: Site ID: Location: ME

Instrument ID: F GC Columns: RTX-CLPesticidesI ID: 0.32(mm)

RTX-CLPesticidesII

Sample ID: 56617-11

COMPOUND	RESPONSE #1 ug/kg	RESPONSE #2 ug/kg	DIFFERENCE %	#	% DIFFERENCE Acceptance Limit
PCB 1260	42	470	167		40

Column to be used to flag % Difference with an asterisk

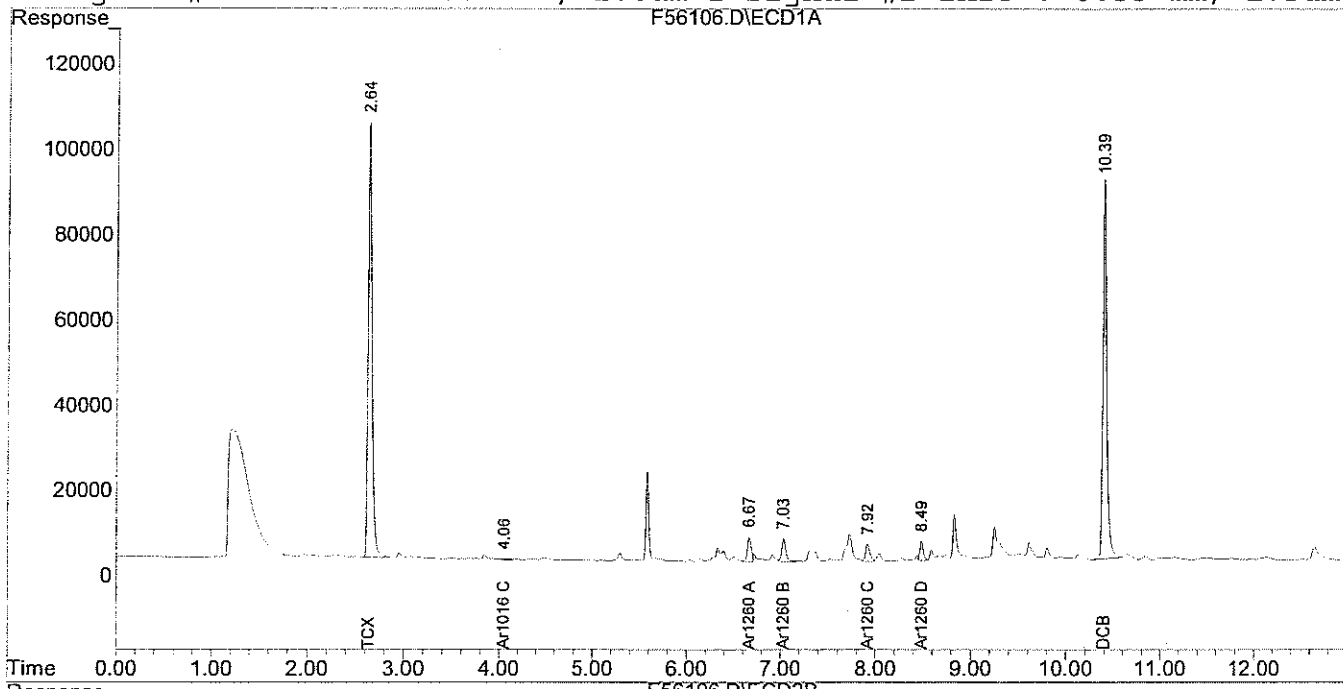
* Values outside of QC limits

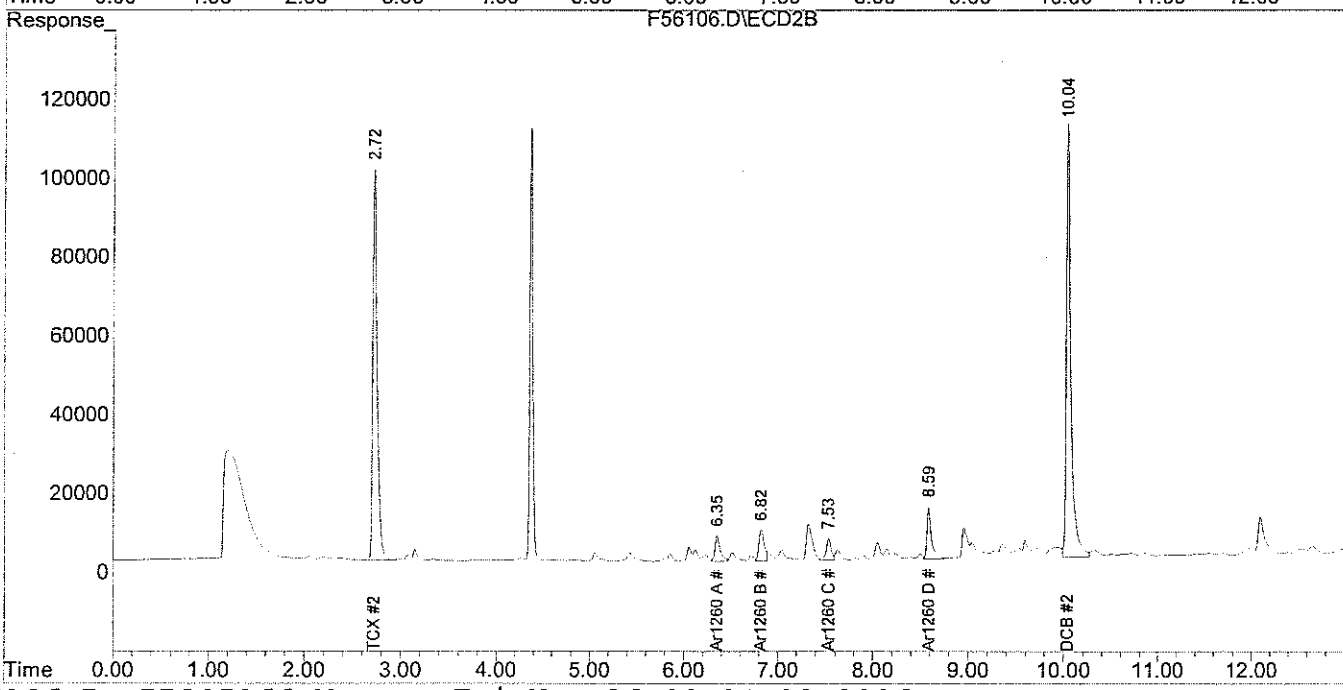
Quantitation Report (QT Reviewed)

Signal #1 : C:\HPCHEM\2\DATA\052506-F\F56106.D\ECD1A.CH Vial: 29
Signal #2 : C:\HPCHEM\2\DATA\052506-F\F56106.D\ECD2B.CH
Acq On : 25 May 2006 10:44 pm Operator:
Sample : 56617-11, A/C Inst : Instrumen
Misc : SOIL Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 26 8:24 2006 Quant Results File: PBQ05166.RES

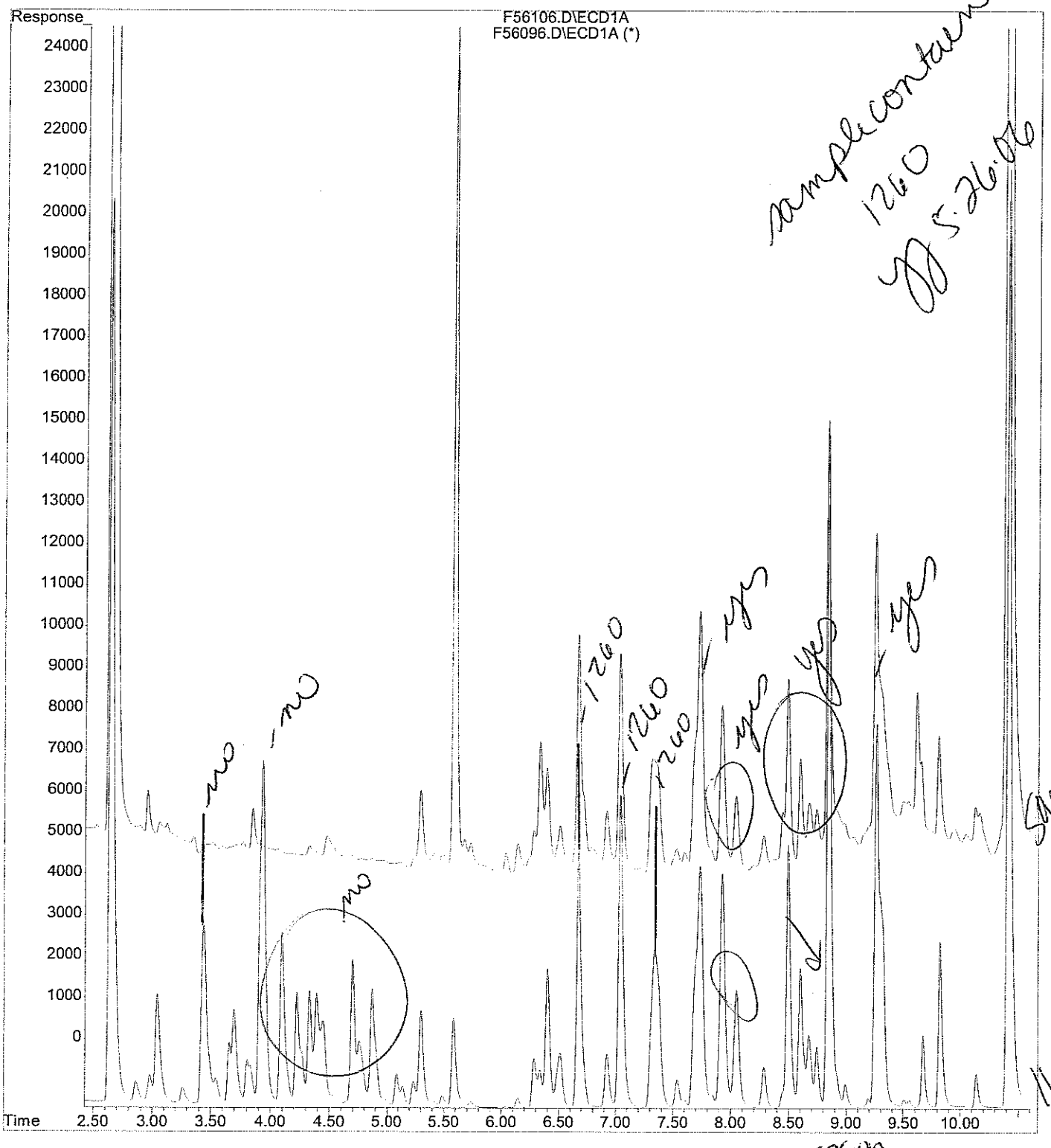
Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 10:06:32 2006
Response via : Multiple Level Calibration
DataAcq Meth : PSTQUICK.M

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film





File : C:\HPCHEM\2\DATA\052506-F\F56106.D
Operator :
Acquired : 25 May 2006 10:44 pm using AcqMethod PSTQUICK.M
Instrument : Instrumen
Sample Name: 56617-11, A/C
Misc Info : SOIL
Vial Number: 29



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May 26, 2006

CLIENT SAMPLE ID

Project Name: GP

Project Number: 218736

Field Sample ID: SC-01

SAMPLE DATA

Lab Sample ID: 56617-12
Matrix: Solid
Percent Solid: 96
Dilution Factor: 1.0
Collection Date: 05/24/06
Lab Receipt Date: 05/25/06
Extraction Date: 05/25/06
Analysis Date: 05/25/06

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit µg/kg	Results µg/kg
PCB-1016	15	U
PCB-1221	15	U
PCB-1232	15	U
PCB-1242	15	U
PCB-1248	15	U
PCB-1254	15	U
PCB-1260	15	U

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	92	%
Decachlorobiphenyl	101	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082.

Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3545.

COMMENTS: Results are expressed on a dry weight basis.

PCB Report

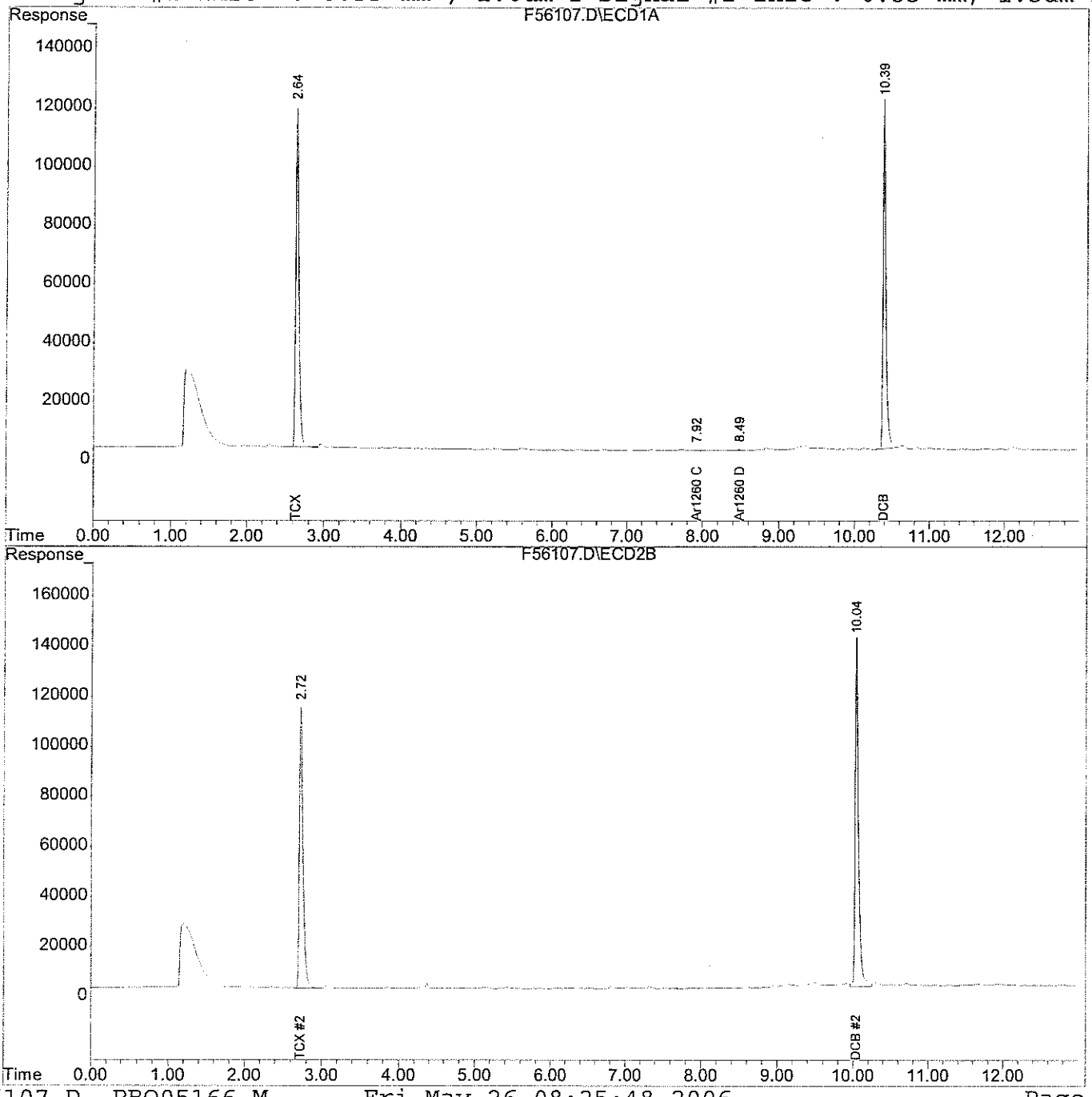
Authorized signature

Quantitation Report (QT Reviewed)

Signal #1 : C:\HPCHEM\2\DATA\052506-F\F56107.D\ECD1A.CH Vial: 30
Signal #2 : C:\HPCHEM\2\DATA\052506-F\F56107.D\ECD2B.CH
Acq On : 25 May 2006 10:58 pm Operator:
Sample : 56617-12, A/C Inst : Instrumen
Misc : SOIL Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 26 8:25 2006 Quant Results File: PBQ05166.RES

Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 10:06:32 2006
Response via : Multiple Level Calibration
DataAcq Meth : PSTQUICK.M

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film



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May 26, 2006

CLIENT SAMPLE ID

Project Name: GP

Project Number: 218736

Field Sample ID: SC-02

SAMPLE DATA

Lab Sample ID: 56617-13
Matrix: Solid
Percent Solid: 96
Dilution Factor: 1.0
Collection Date: 05/24/06
Lab Receipt Date: 05/25/06
Extraction Date: 05/25/06
Analysis Date: 05/25/06

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit µg/kg	Results µg/kg
PCB-1016	15	U
PCB-1221	15	U
PCB-1232	15	U
PCB-1242	15	**99**
PCB-1248	15	U
PCB-1254	15	U
PCB-1260	15	U

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	92	%
Decachlorobiphenyl	100	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082.

Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3545.

COMMENTS: Results are expressed on a dry weight basis.

PCB Report

Authorized signature Melissa Hall

analytics environmental laboratory LLC

PCB CONFIRMATION RESULTS

SDG: 56617

Project No: Site ID: Location: ME

Instrument ID: F GC Columns: RTX-CLPesticidesI ID: 0.32(mm)

RTX-CLPesticidesII

Sample ID: 56617-13

COMPOUND	RESPONSE #1 ug/kg	RESPONSE #2 ug/kg	DIFFERENCE %	#	% DIFFERENCE Acceptance Limit
PCB 1242	99	108	9		40

Column to be used to flag % Difference with an asterisk

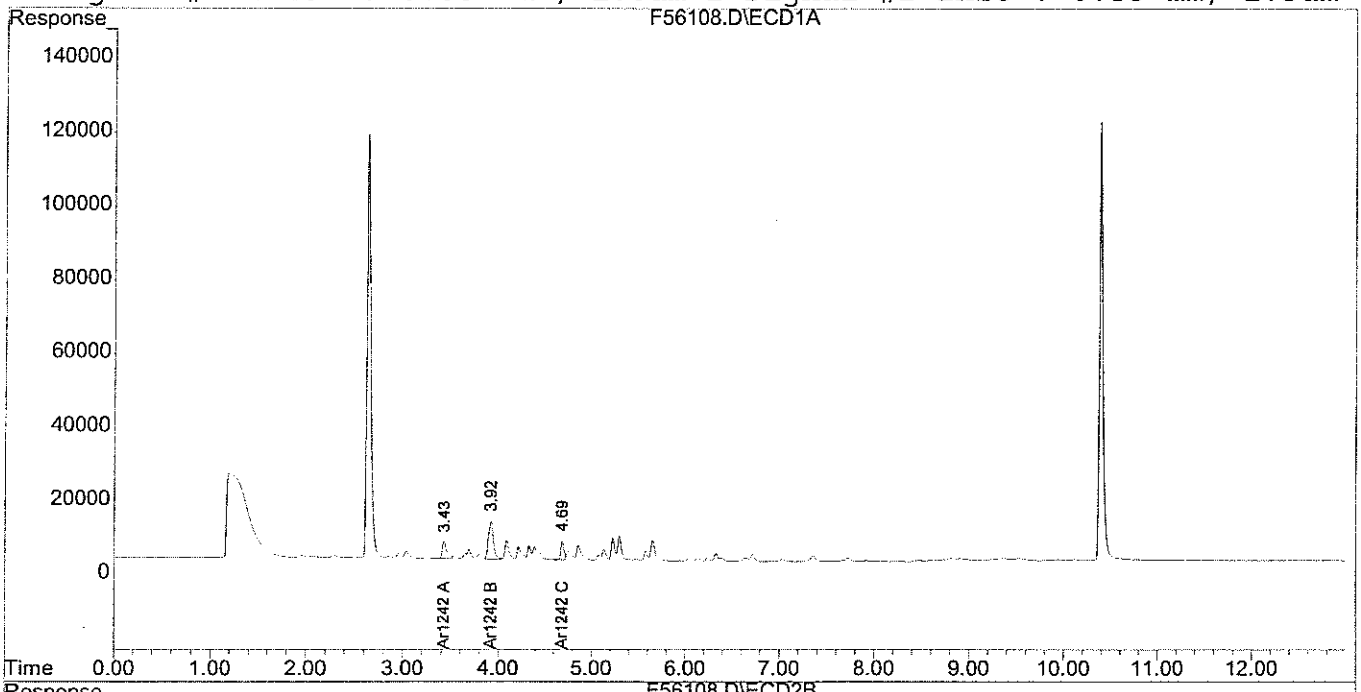
* Values outside of QC limits

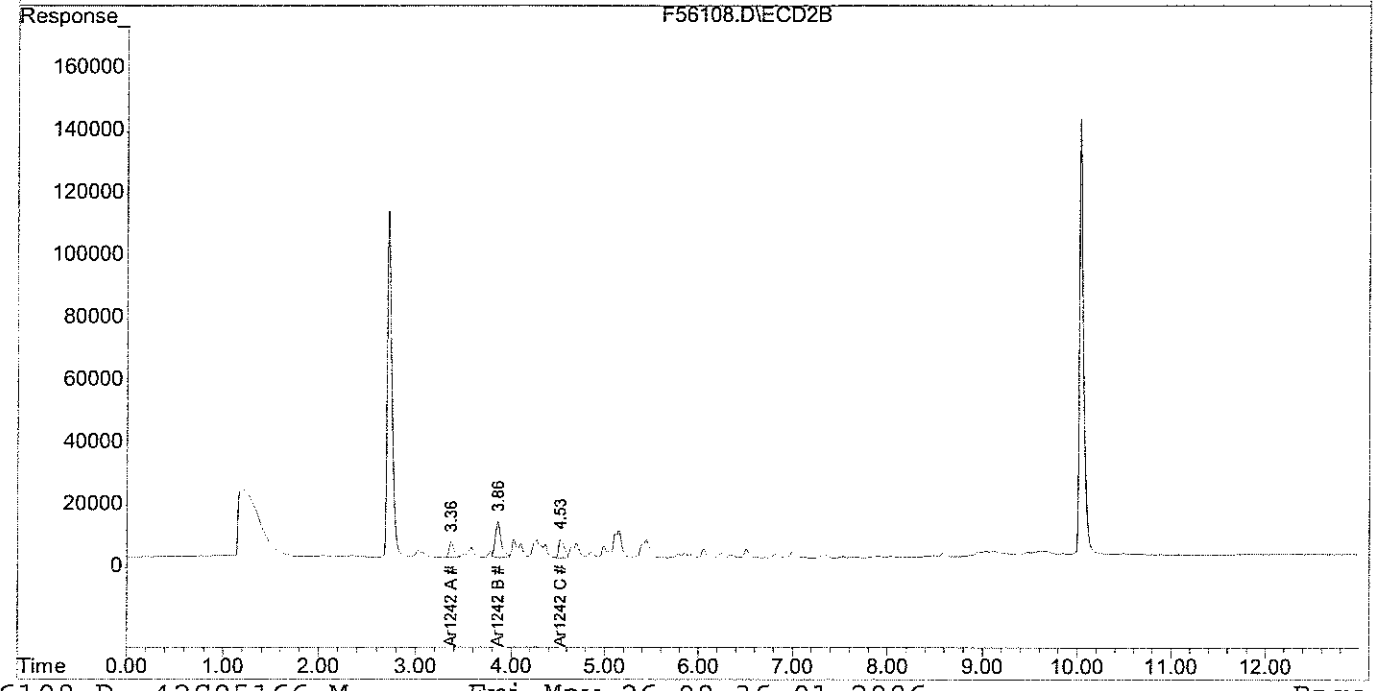
Quantitation Report (QT Reviewed)

Signal #1 : C:\HPCHEM\2\DATA\052506-F\F56108.D\ECD1A.CH Vial: 31
Signal #2 : C:\HPCHEM\2\DATA\052506-F\F56108.D\ECD2B.CH
Acq On : 25 May 2006 11:13 pm Operator:
Sample : 56617-13, A/C Inst : Instrumen
Misc : SOIL Multiplr: 1.00
IntFile Signal #1: EVENTS.E IntFile Signal #2: EVENTS2.E
Quant Time: May 26 8:35 2006 Quant Results File: 42S05166.RES

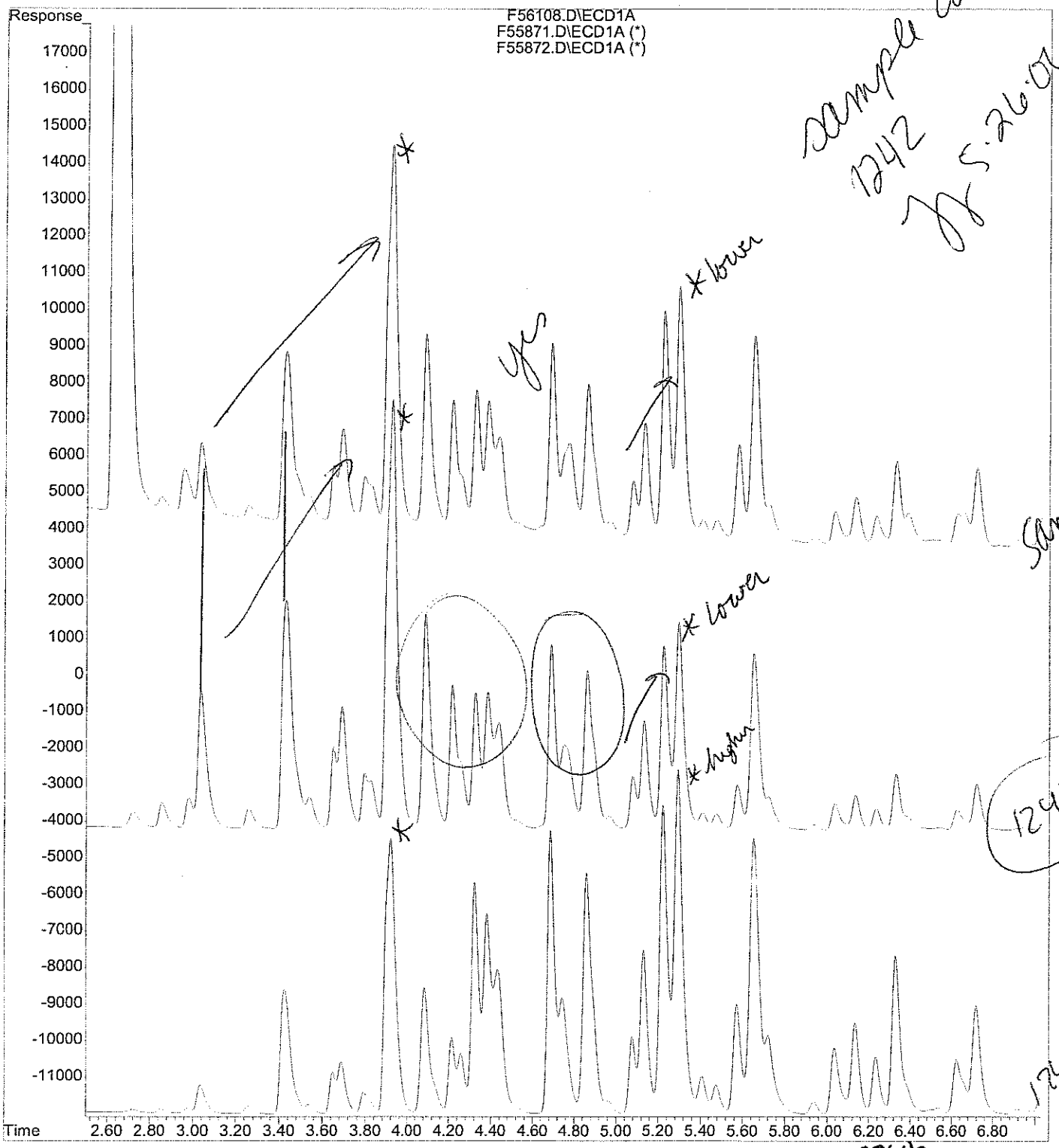
Quant Method : C:\HPCHEM\2\METHODS\42S05166.M (Chemstation Integrator)
Title : Aroclor 1242
Last Update : Fri May 26 08:34:45 2006
Response via : Single Level Calibration
DataAcq Meth : PSTQUICK.M

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film





File : C:\HPCHEM\2\DATA\052506-F\F56108.D
Operator :
Acquired : 25 May 2006 11:13 pm using AcqMethod PSTQUICK.M
Instrument : Instrumen
Sample Name: 56617-13, A/C
Misc Info : SOIL
Vial Number: 31



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May 30, 2006

CLIENT SAMPLE ID

Project Name: GP

Project Number: 218736

Field Sample ID: SC-03

SAMPLE DATA

Lab Sample ID: 56617-14 DL
Matrix: Solid
Percent Solid: 98
Dilution Factor: 6910
Collection Date: 05/24/06
Lab Receipt Date: 05/25/06
Extraction Date: 05/25/06
Analysis Date: 05/26/06

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit µg/kg	Results µg/kg
PCB-1016	104000	U
PCB-1221	104000	U
PCB-1232	104000	U
PCB-1242	104000	U
PCB-1248	104000	U
PCB-1254	104000	U
PCB-1260	104000	**4300000**

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	*	%
Decachlorobiphenyl	*	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082.

Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3545.

COMMENTS: Results are expressed on a dry weight basis. * The surrogates were diluted out.

PCB Report

Authorized signature Melissa Hull

analytics environmental laboratory LLC

PCB CONFIRMATION RESULTS

SDG: 56617

Project No: Site ID: Location: ME

Instrument ID: F GC Columns: RTX-CLPesticidesI ID: 0.32(mm)

RTX-CLPesticidesII

Sample ID: 56617-14

COMPOUND	RESPONSE #1 ug/kg	RESPONSE #2 ug/kg	DIFFERENCE %	#	% DIFFERENCE Acceptance Limit
PCB 1260	4300000	4330000	1		40

Column to be used to flag % Difference with an asterisk

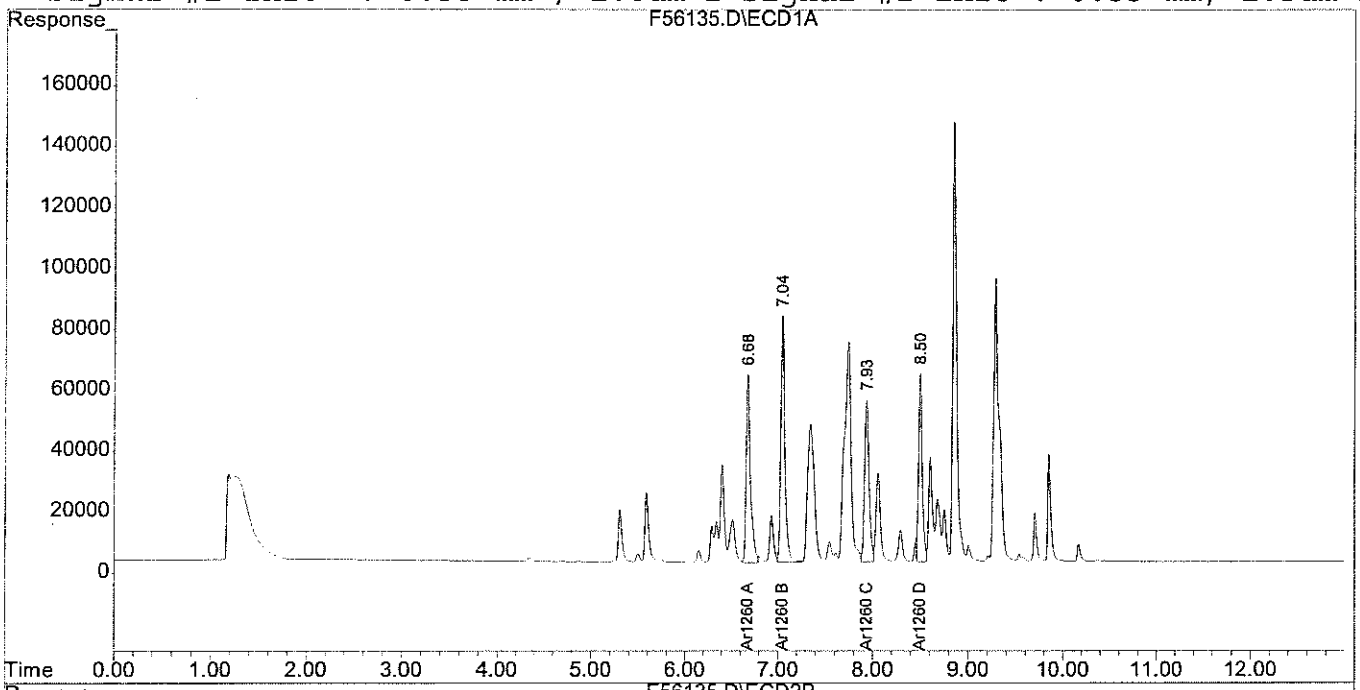
* Values outside of QC limits

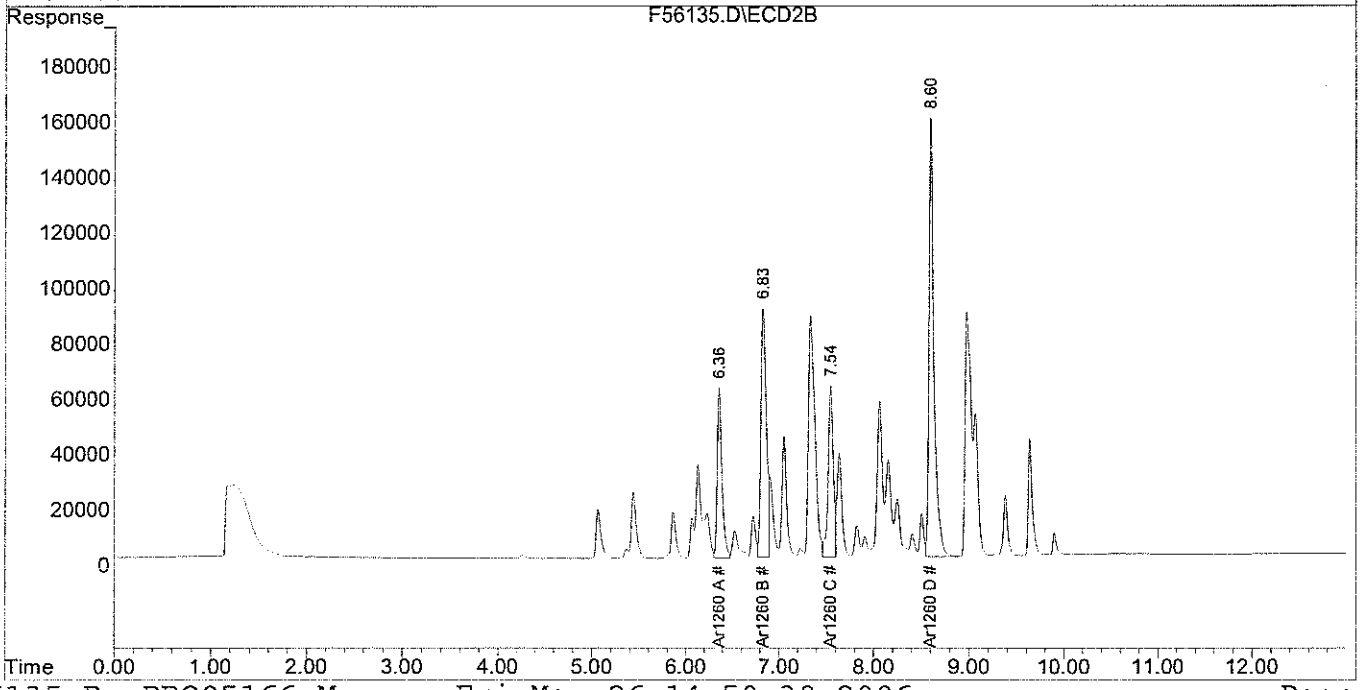
Quantitation Report (QT Reviewed)

Signal #1 : C:\HPCHEM\2\DATA\052606-F\F56135.D\ECD1A.CH Vial: 52
Signal #2 : C:\HPCHEM\2\DATA\052606-F\F56135.D\ECD2B.CH
Acq On : 26 May 2006 2:36 pm Operator:
Sample : 56617-14, 1:2000, A/C Inst : Instrumen
Misc : SOIL Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 26 14:50 2006 Quant Results File: PBQ05166.RES

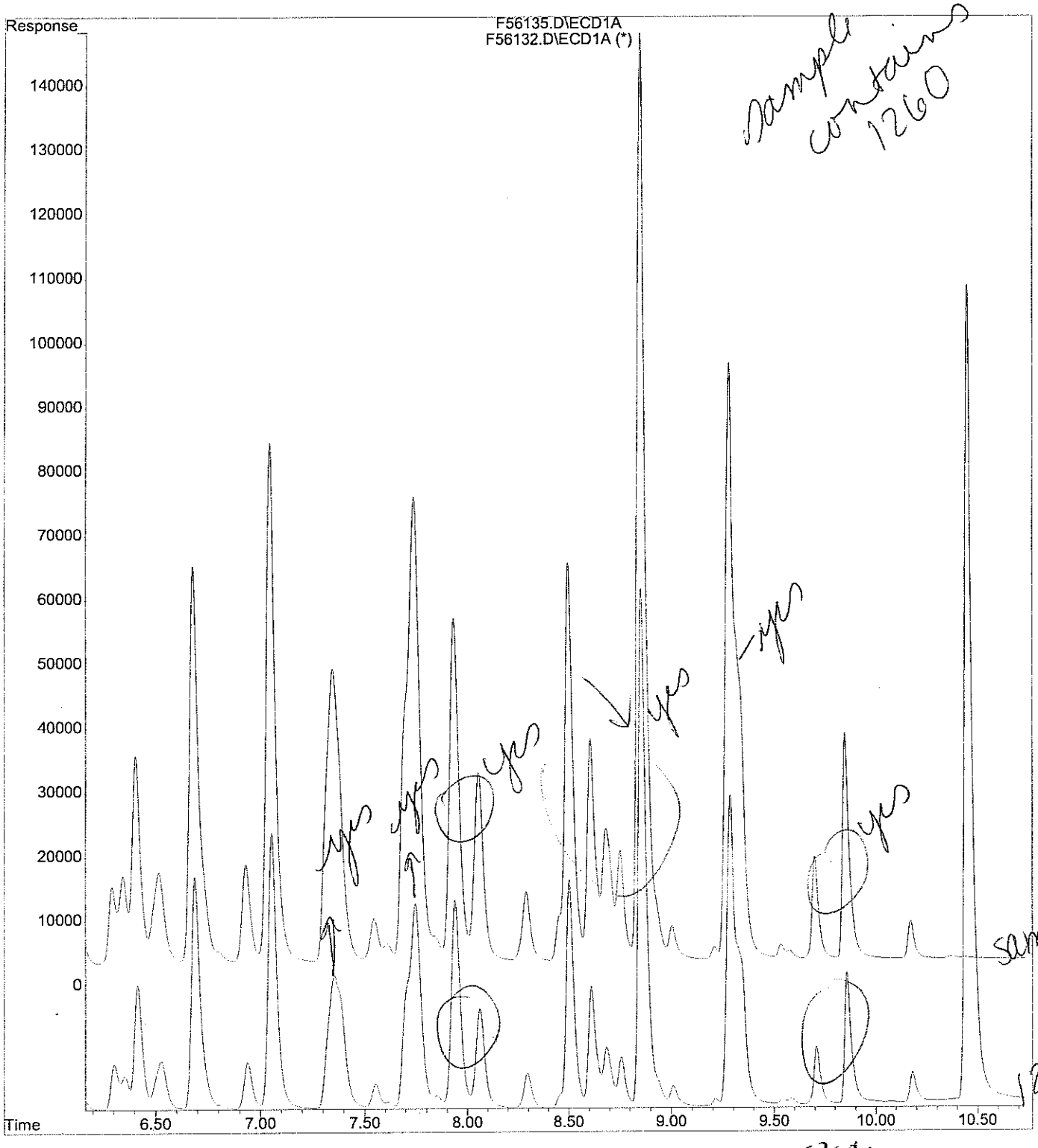
Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 10:06:32 2006
Response via : Multiple Level Calibration
DataAcq Meth : PSTQUICK.M

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film





File : C:\HPCHEM\2\DATA\052606-F\F56135.D
Operator :
Acquired : 26 May 2006 2:36 pm using AcqMethod PSTQUICK.M
Instrument : Instrumen
Sample Name: 56617-14, 1:2000, A/C
Misc Info : SOIL
Vial Number: 52



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Mr. Karl Kasper
Woodard & Curran
41 Hutchins Drive
Portland ME 04102

May 26, 2006

SAMPLE DATA

CLIENT SAMPLE ID

Project Name:	GP
Project Number:	218736
Field Sample ID:	SS-05

Lab Sample ID:	56617-15
Matrix:	Solid
Percent Solid:	83
Dilution Factor:	1.1
Collection Date:	05/24/06
Lab Receipt Date:	05/25/06
Extraction Date:	05/26/06
Analysis Date:	05/26/06

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit µg/kg	Results µg/kg
PCB-1016	17	U
PCB-1221	17	U
PCB-1232	17	U
PCB-1242	17	U
PCB-1248	17	U
PCB-1254	17	U
PCB-1260	17	U

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	85	%
Decachlorobiphenyl	75	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082.

Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3545.

COMMENTS: Results are expressed on a dry weight basis.

PCB Report

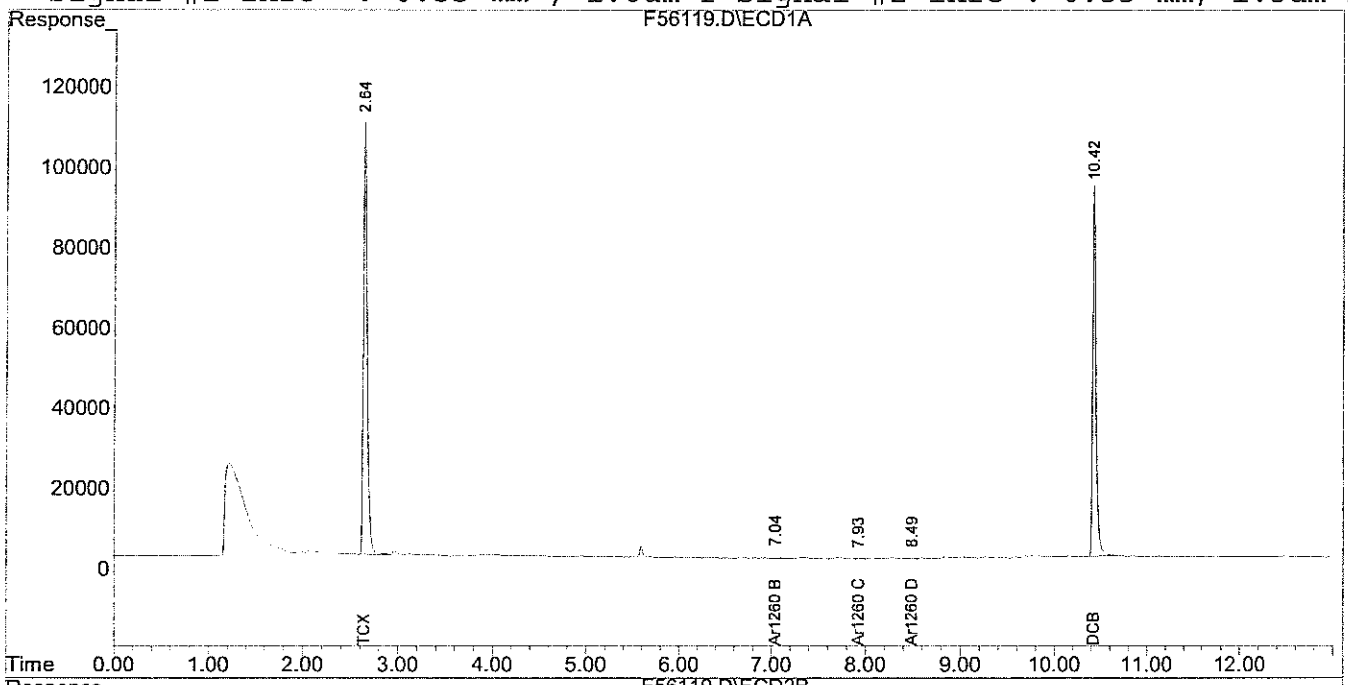
Authorized signature Melissa Hall

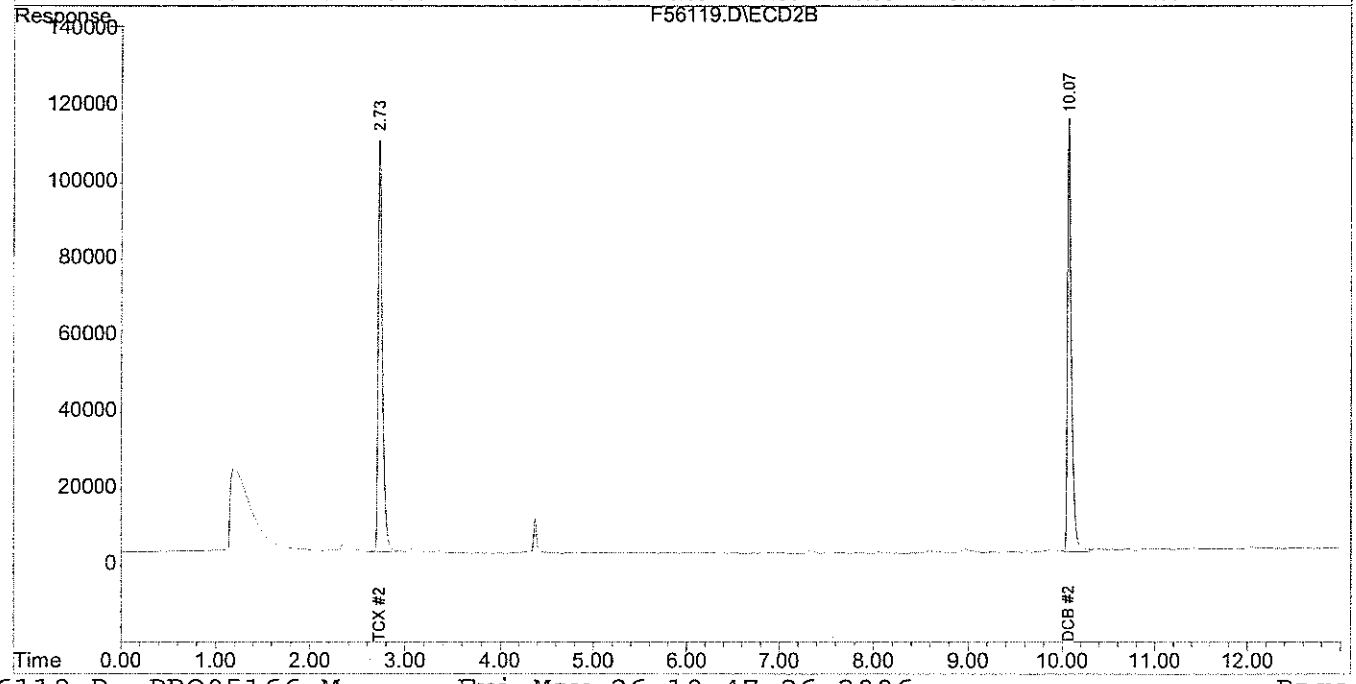
Quantitation Report (Not Reviewed)

Signal #1 : C:\HPCHEM\2\DATA\052506-F\F56119.D\ECD1A.CH Vial: 39
Signal #2 : C:\HPCHEM\2\DATA\052506-F\F56119.D\ECD2B.CH
Acq On : 26 May 2006 8:18 am Operator:
Sample : 56617-15, A/C Inst : Instrumen
Misc : SOIL Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 26 10:47 2006 Quant Results File: PBQ05166.RES

Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 10:06:32 2006
Response via : Multiple Level Calibration
DataAcq Meth : PSTQUICK.M

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film





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Portland ME 04102

May 30, 2006

CLIENT SAMPLE ID

Project Name: GP

Project Number: 218736

Field Sample ID: SS-06

SAMPLE DATA

Lab Sample ID: 56617-16 DL
Matrix: Solid
Percent Solid: 93
Dilution Factor: 10
Collection Date: 05/24/06
Lab Receipt Date: 05/25/06
Extraction Date: 05/26/06
Analysis Date: 05/26/06

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit µg/kg	Results µg/kg
PCB-1016	150	U
PCB-1221	150	U
PCB-1232	150	U
PCB-1242	150	U
PCB-1248	150	U
PCB-1254	150	**2720**
PCB-1260	150	**3330**

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	90	%
Decachlorobiphenyl	120	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082.

Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3545.

COMMENTS: Results are expressed on a dry weight basis.

PCB Report

Authorized signature Melissa Hall

analytics environmental laboratory LLC

PCB CONFIRMATION RESULTS

SDG: 56617

Project No: Site ID: Location: ME

Instrument ID: F GC Columns: RTX-CLPesticidesI ID: 0.32(mm)
RTX-CLPesticidesII

Sample ID: 56617-16

COMPOUND	RESPONSE #1 ug/kg	RESPONSE #2 ug/kg	DIFFERENCE %	#	% DIFFERENCE Acceptance Limit
PCB 1254	2720	2550	6		40
PCB 1260	3330	3470	4		40

Column to be used to flag % Difference with an asterisk

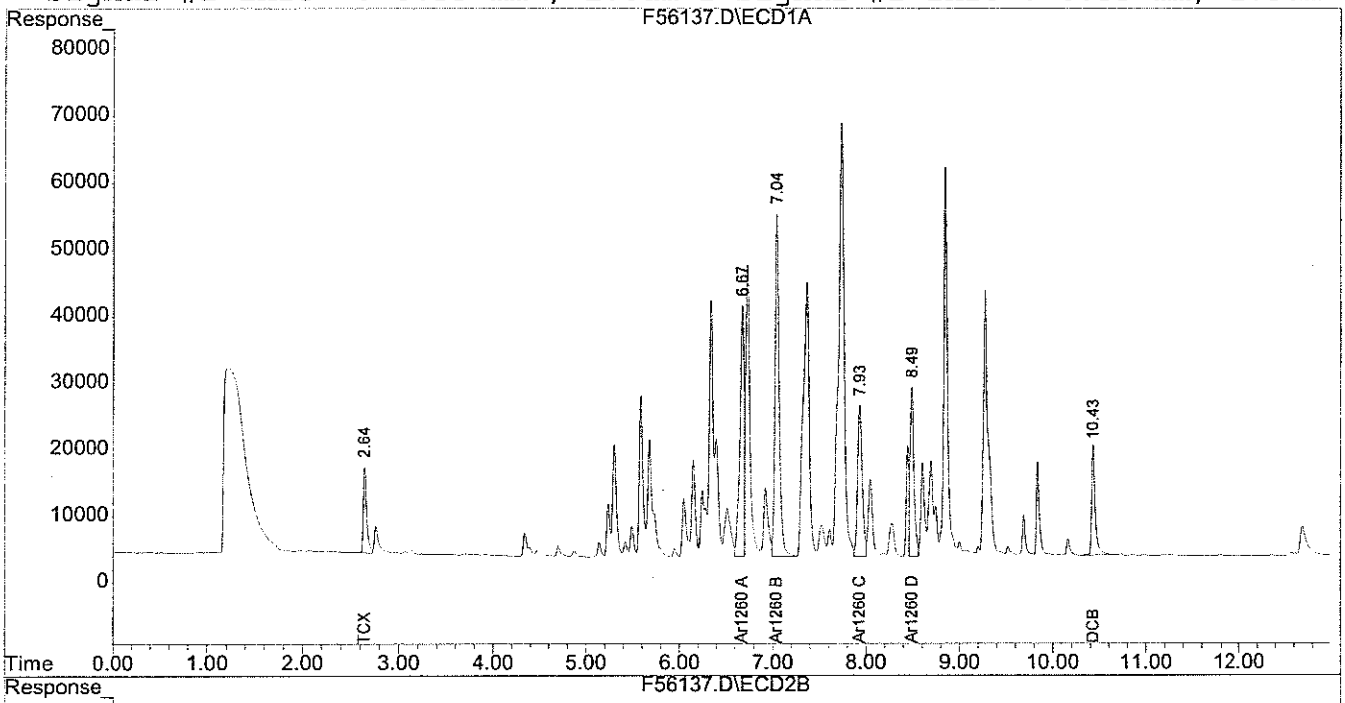
* Values outside of QC limits

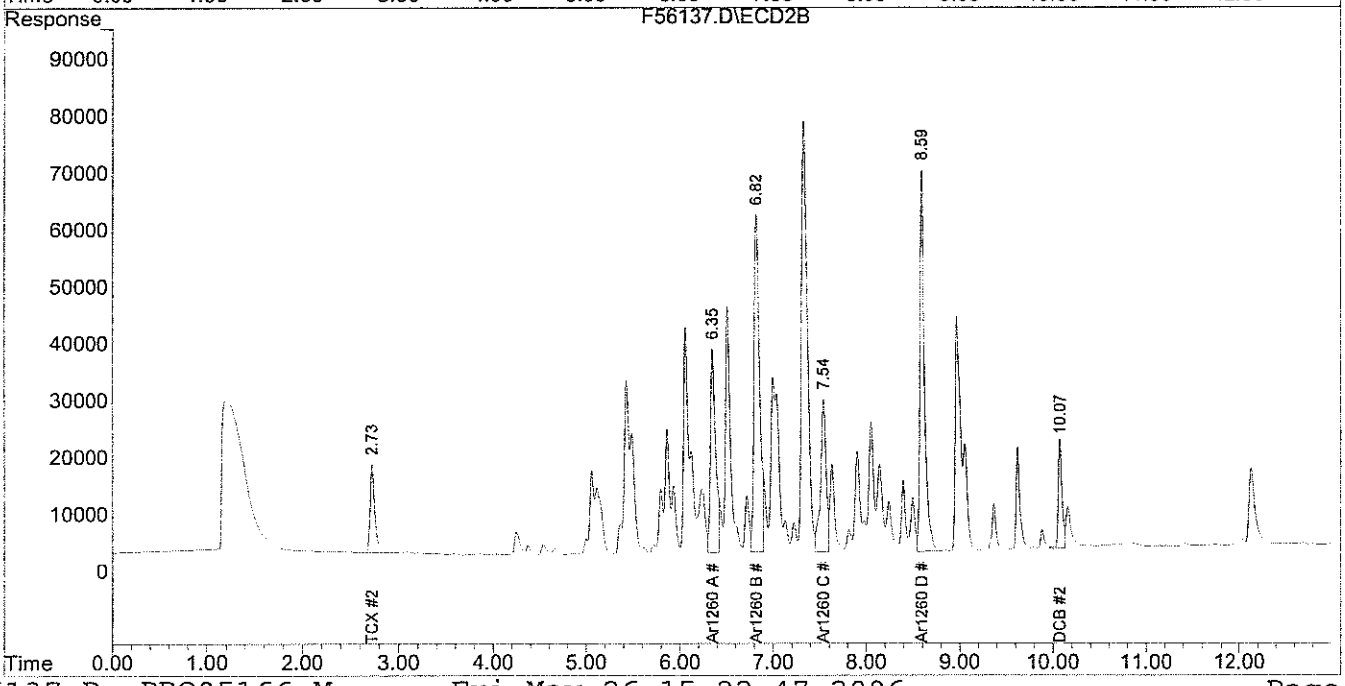
Quantitation Report (QT Reviewed)

Signal #1 : C:\HPCHEM\2\DATA\052606-F\F56137.D\ECD1A.CH Vial: 54
Signal #2 : C:\HPCHEM\2\DATA\052606-F\F56137.D\ECD2B.CH
Acq On : 26 May 2006 3:05 pm Operator:
Sample : 56617-16, 1:10, A/C Inst : Instrumen
Misc : SOIL Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 26 15:22 2006 Quant Results File: PBQ05166.RES

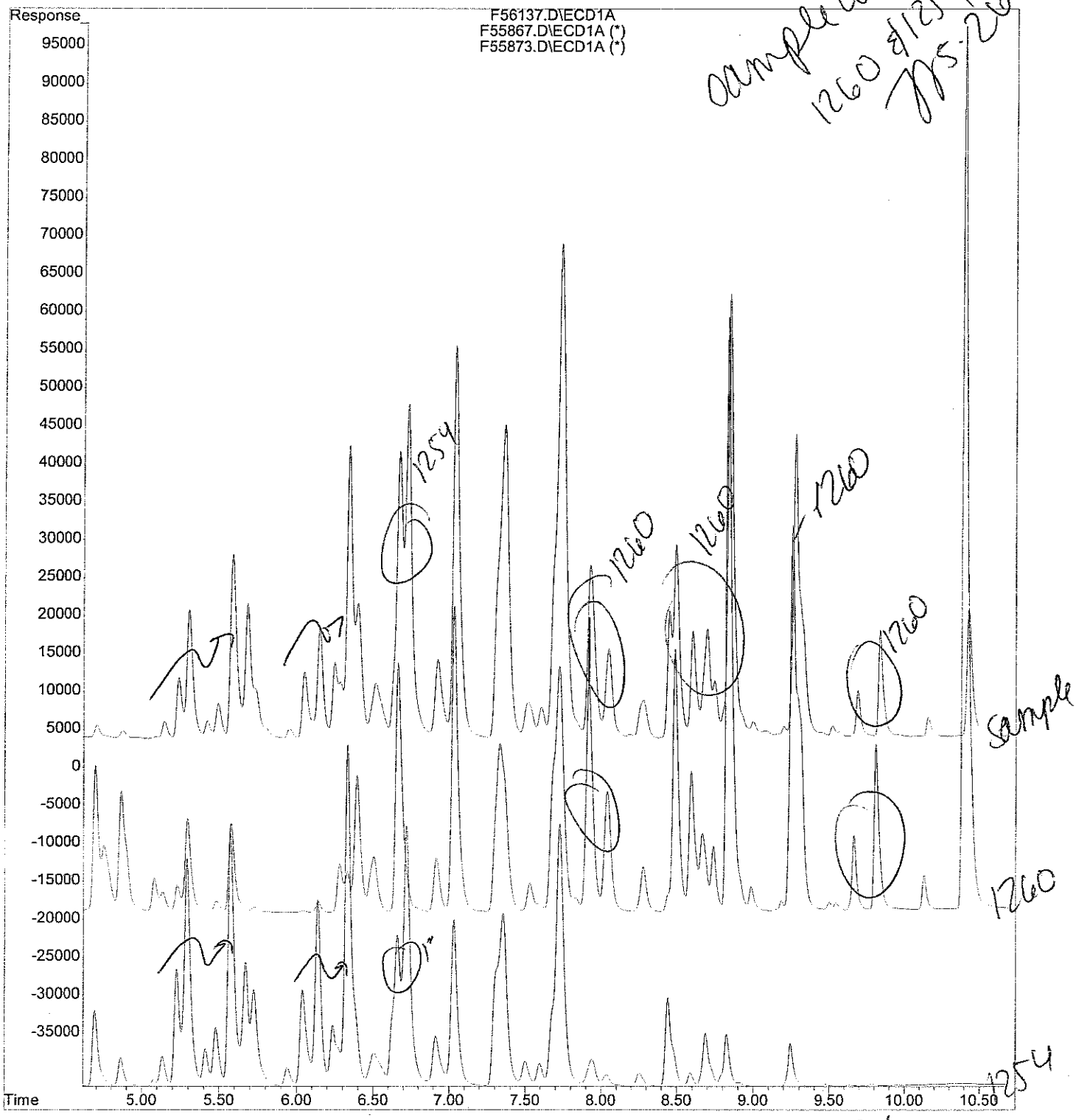
Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 10:06:32 2006
Response via : Multiple Level Calibration
DataAcq Meth : PSTQUICK.M

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film





File : C:\HPCHEM\2\DATA\052606-F\F56137.D
Operator :
Acquired : 26 May 2006 3:05 pm using AcqMethod PSTQUICK.M
Instrument : Instrumen
Sample Name: 56617-16, 1:10, A/C
Misc Info : SOIL
Vial Number: 54



sample conditions 1260 & 1254 5-26-06

5-26-06

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Portland ME 04102

May 30, 2006

CLIENT SAMPLE ID

Project Name: GP

Project Number: 218736

Field Sample ID: SS-07

SAMPLE DATA

Lab Sample ID: 56617-17 DL
Matrix: Solid
Percent Solid: 95
Dilution Factor: 52
Collection Date: 05/24/06
Lab Receipt Date: 05/25/06
Extraction Date: 05/26/06
Analysis Date: 05/26/06

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit µg/kg	Results µg/kg
PCB-1016	780	U
PCB-1221	780	U
PCB-1232	780	U
PCB-1242	780	U
PCB-1248	780	U
PCB-1254	780	**7650**
PCB-1260	780	**15500**

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	*	%
Decachlorobiphenyl	*	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082.

COMMENTS: Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3545. Results are expressed on a dry weight basis. * The surrogates were diluted out.

PCB Report

Authorized signature Melinda Hall

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PCB CONFIRMATION RESULTS

SDG: 56617

Project No: Site ID: Location: ME

Instrument ID: F GC Columns: RTX-CLPesticidesI ID: 0.32(mm)
RTX-CLPesticidesII

Sample ID: 56617-17

COMPOUND	RESPONSE #1 ug/kg	RESPONSE #2 ug/kg	DIFFERENCE %	#	% DIFFERENCE Acceptance Limit
PCB 1254	7650	7580	1		40
PCB 1260	15500	17000	9		40

Column to be used to flag % Difference with an asterisk

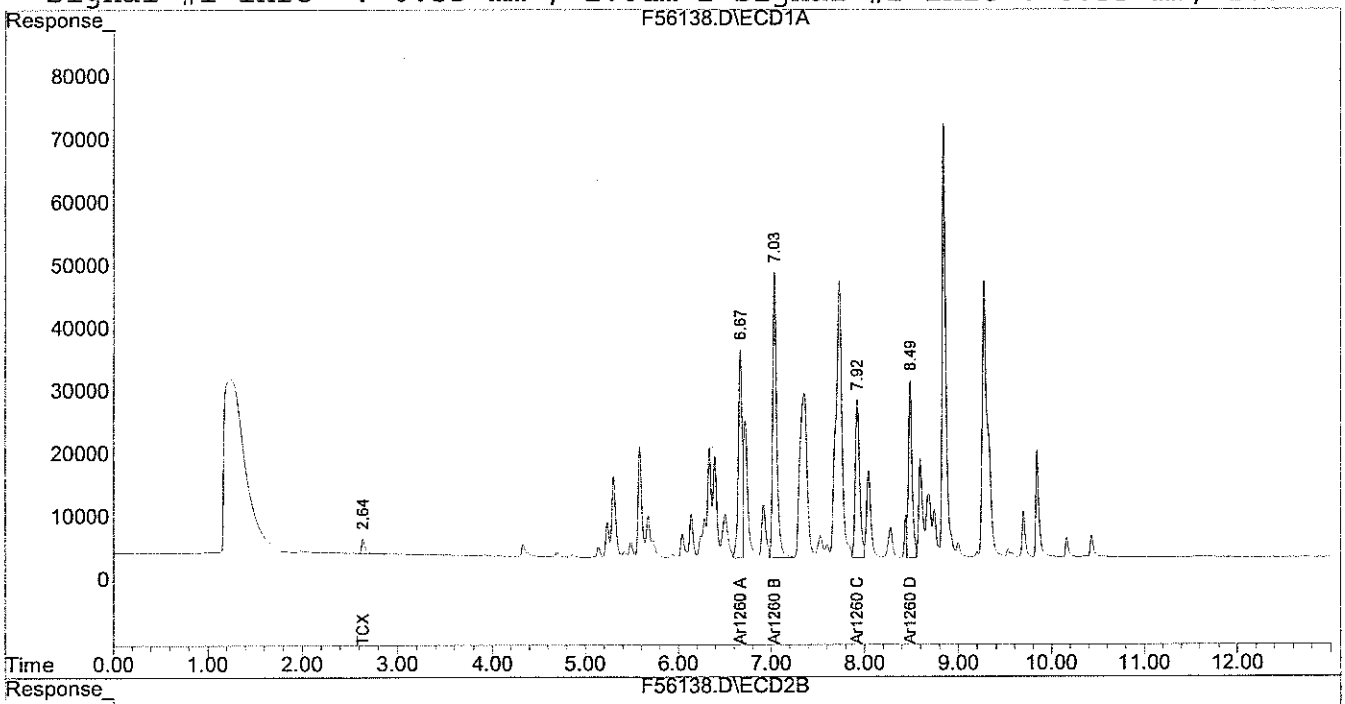
* Values outside of QC limits

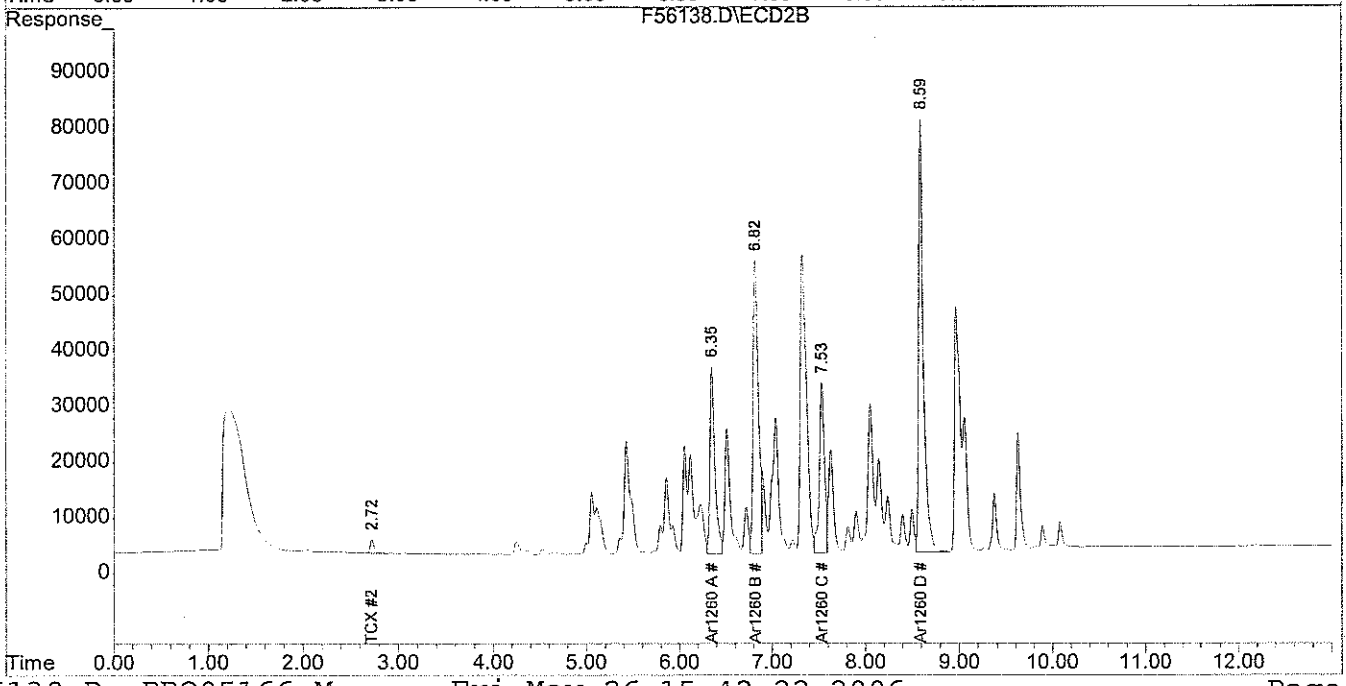
Quantitation Report (QT Reviewed)

Signal #1 : C:\HPCHEM\2\DATA\052606-F\F56138.D\ECD1A.CH Vial: 55
Signal #2 : C:\HPCHEM\2\DATA\052606-F\F56138.D\ECD2B.CH
Acq On : 26 May 2006 3:35 pm Operator:
Sample : 56617-17, 1:50, A/C Inst : Instrumen
Misc : SOIL Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 26 15:42 2006 Quant Results File: PBQ05166.RES

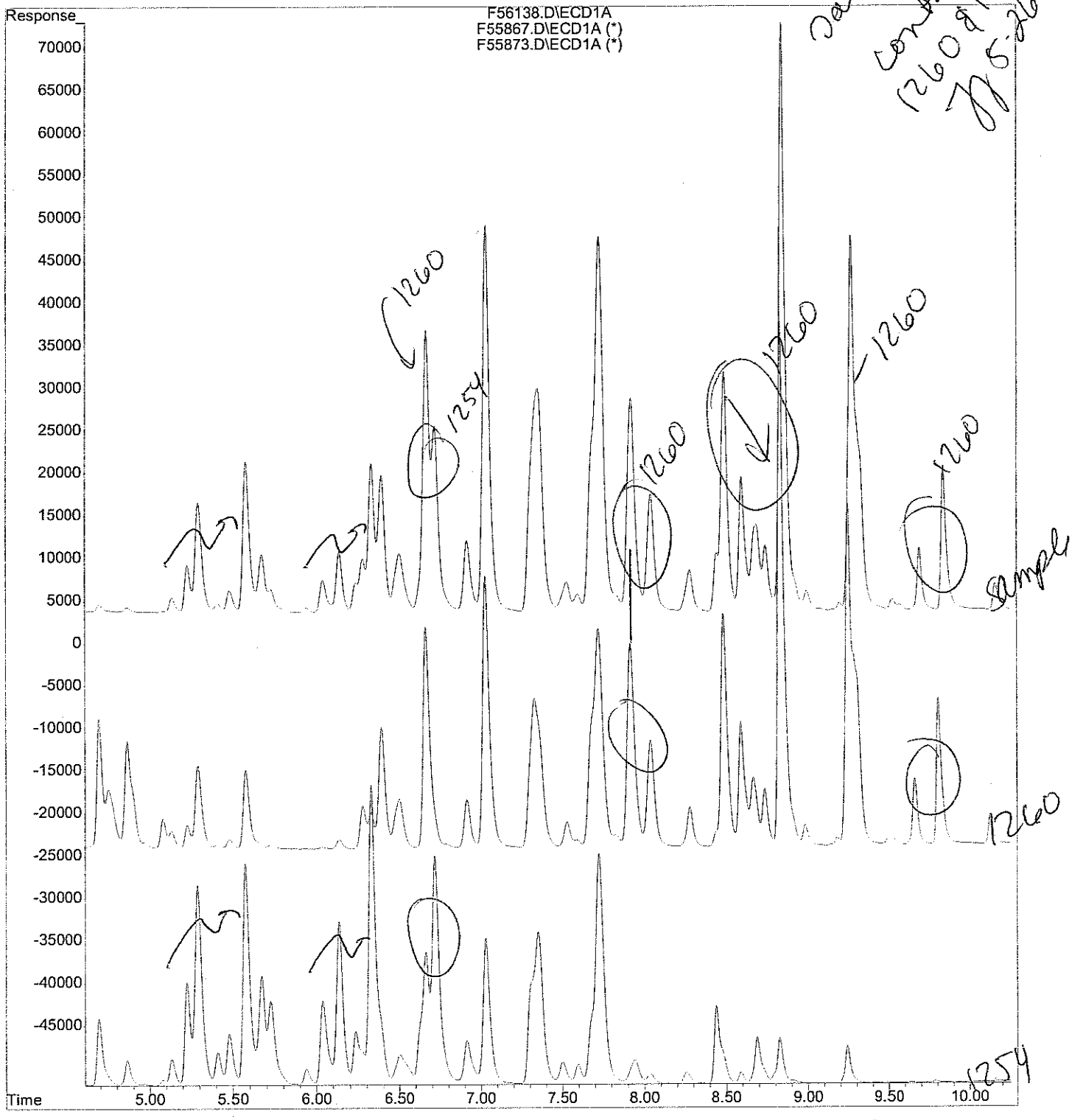
Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 10:06:32 2006
Response via : Multiple Level Calibration
DataAcq Meth : PSTQUICK.M

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film





File : C:\HPCHEM\2\DATA\052606-F\F56138.D
Operator :
Acquired : 26 May 2006 3:35 pm using AcqMethod PSTQUICK.M
Instrument : Instrumen
Sample Name: 56617-17, 1:50, A/C
Misc Info : SOIL
Vial Number: 55



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Portland ME 04102

May 26, 2006

CLIENT SAMPLE ID

Project Name: GP

Project Number: 218736

Field Sample ID: SS-08

SAMPLE DATA

Lab Sample ID: 56617-18
Matrix: Solid
Percent Solid: 94
Dilution Factor: 1.1
Collection Date: 05/24/06
Lab Receipt Date: 05/25/06
Extraction Date: 05/26/06
Analysis Date: 05/26/06

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit µg/kg	Results µg/kg
PCB-1016	17	U
PCB-1221	17	U
PCB-1232	17	U
PCB-1242	17	U
PCB-1248	17	U
PCB-1254	17	U
PCB-1260	17	U

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	102	%
Decachlorobiphenyl	97	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082.

Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3545.

COMMENTS: Results are expressed on a dry weight basis.

PCB Report

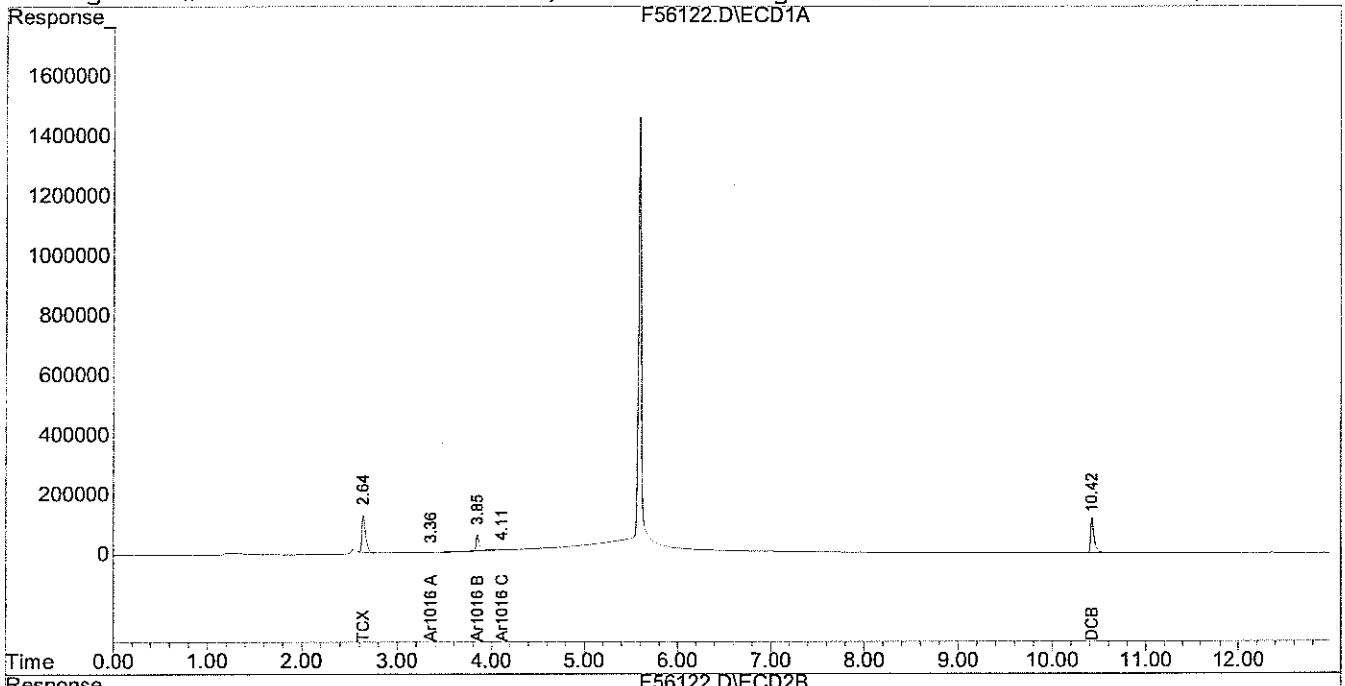
Authorized signature Melissa Hall

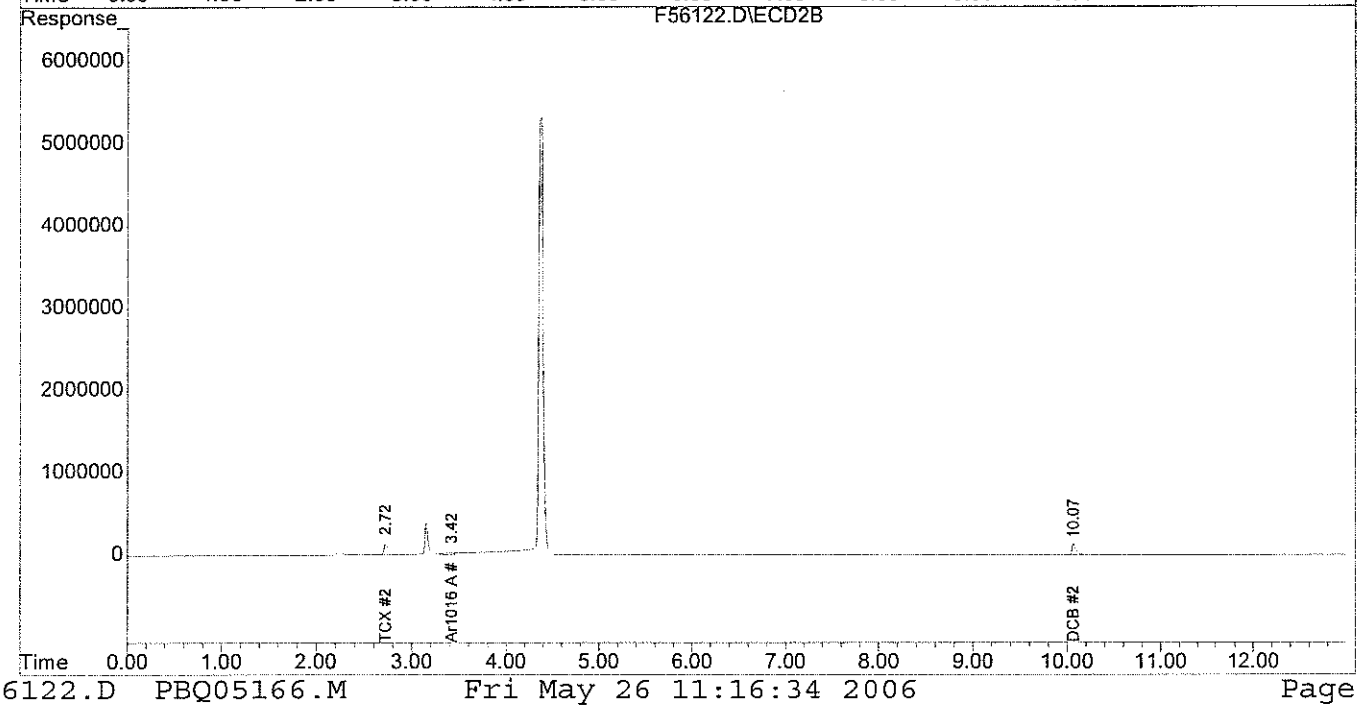
Quantitation Report (QT Reviewed)

Signal #1 : C:\HPCHEM\2\DATA\052506-F\F56122.D\ECD1A.CH Vial: 42
Signal #2 : C:\HPCHEM\2\DATA\052506-F\F56122.D\ECD2B.CH
Acq On : 26 May 2006 9:01 am Operator:
Sample : 56617-18, A/C Inst : Instrumen
Misc : SOIL Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 26 11:16 2006 Quant Results File: PBQ05166.RES

Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 10:06:32 2006
Response via : Multiple Level Calibration
DataAcq Meth : PSTQUICK.M

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film





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Portland ME 04102

May 26, 2006

CLIENT SAMPLE ID

Project Name: GP

Project Number: 218736

Field Sample ID: SS-09

SAMPLE DATA

Lab Sample ID: 56617-19
Matrix: Solid
Percent Solid: 94
Dilution Factor: 1.0
Collection Date: 05/24/06
Lab Receipt Date: 05/25/06
Extraction Date: 05/26/06
Analysis Date: 05/26/06

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit µg/kg	Results µg/kg
PCB-1016	15	U
PCB-1221	15	U
PCB-1232	15	U
PCB-1242	15	U
PCB-1248	15	U
PCB-1254	15	U
PCB-1260	15	U

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	96	%
Decachlorobiphenyl	95	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082.

Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3545.

COMMENTS: Results are expressed on a dry weight basis.

PCB Report

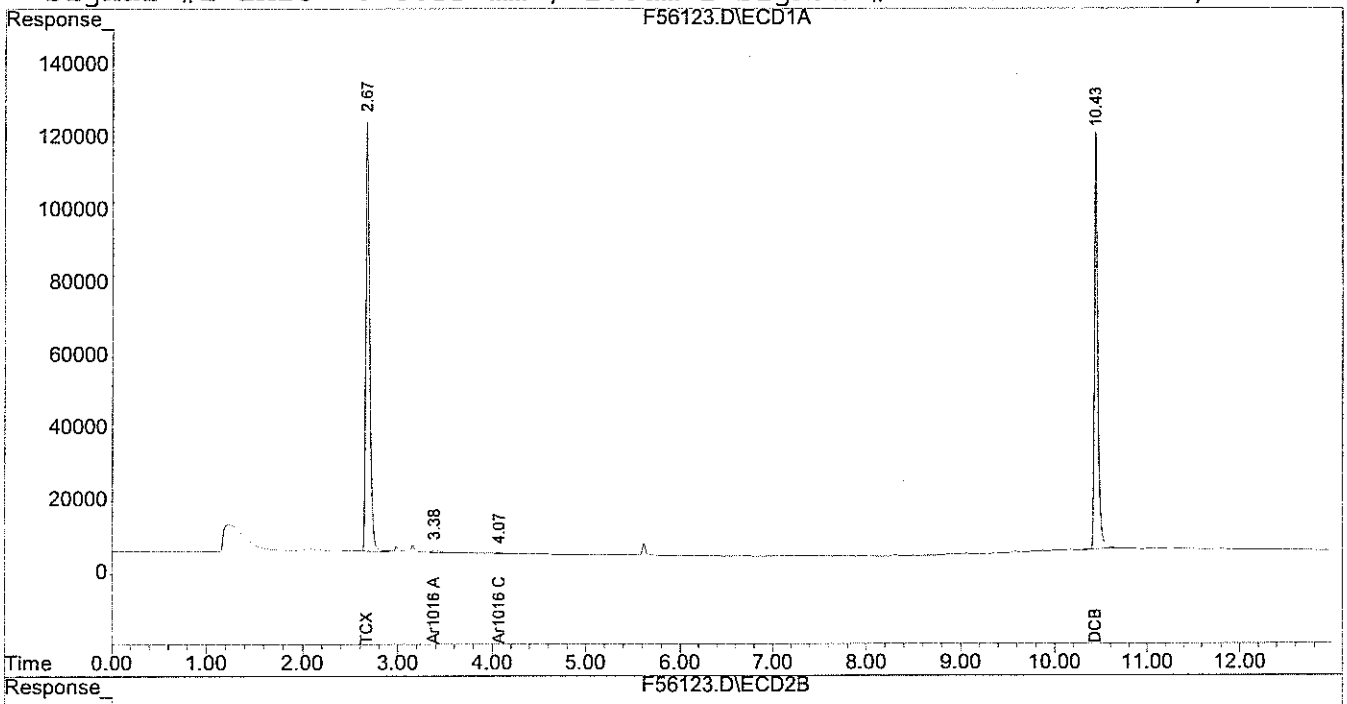
Authorized signature

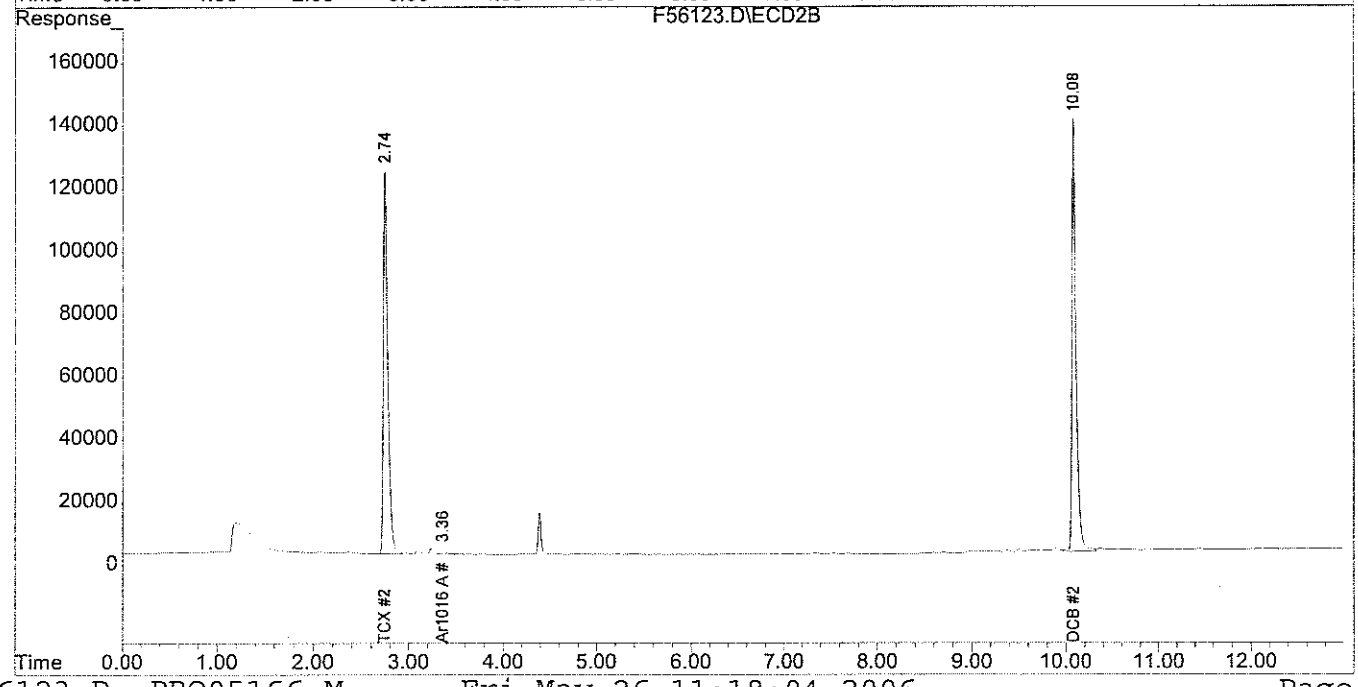
Quantitation Report (QT Reviewed)

Signal #1 : C:\HPCHEM\2\DATA\052506-F\F56123.D\ECD1A.CH Vial: 43
Signal #2 : C:\HPCHEM\2\DATA\052506-F\F56123.D\ECD2B.CH
Acq On : 26 May 2006 9:31 am Operator:
Sample : 56617-19, A/C Inst : Instrumen
Misc : SOIL Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 26 11:17 2006 Quant Results File: PBQ05166.RES

Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 10:06:32 2006
Response via : Multiple Level Calibration
DataAcq Meth : PSTQUICK.M

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film





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Portland ME 04102

May 26, 2006

CLIENT SAMPLE ID

Project Name: GP

Project Number: 218736

Field Sample ID: SS-09 DUP

SAMPLE DATA

Lab Sample ID: 56617-20
Matrix: Solid
Percent Solid: 90
Dilution Factor: 1.1
Collection Date: 05/24/06
Lab Receipt Date: 05/25/06
Extraction Date: 05/26/06
Analysis Date: 05/26/06

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit µg/kg	Results µg/kg
PCB-1016	17	U
PCB-1221	17	U
PCB-1232	17	U
PCB-1242	17	U
PCB-1248	17	U
PCB-1254	17	U
PCB-1260	17	U

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	100	%
Decachlorobiphenyl	95	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082.

Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3545.

COMMENTS: Results are expressed on a dry weight basis.

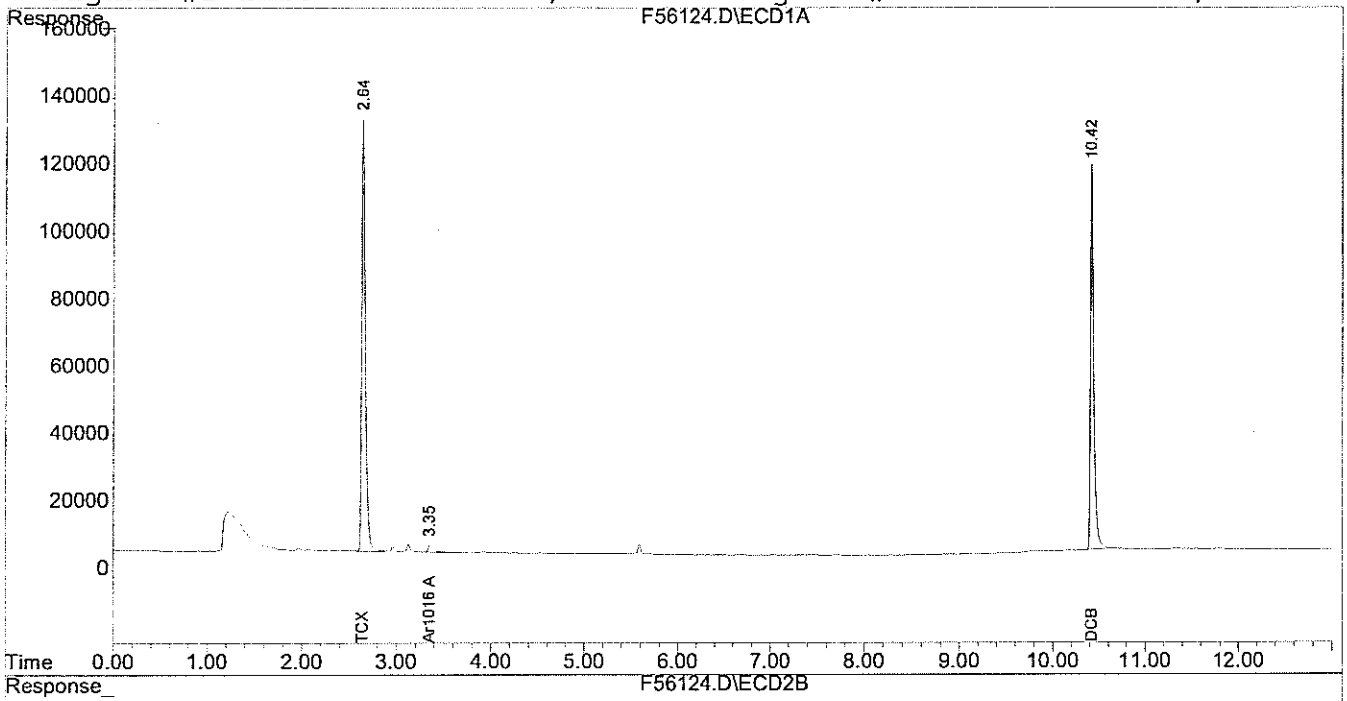
PCB Report

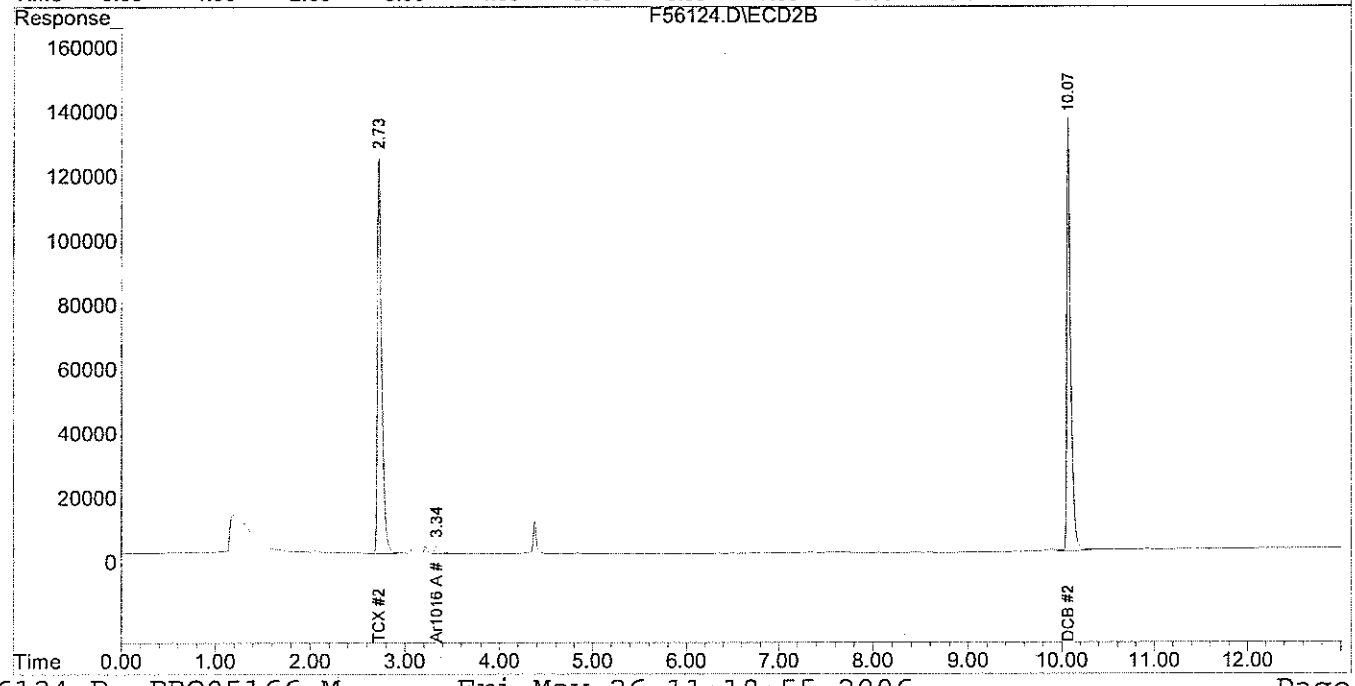
Authorized signature

Signal #1 : C:\HPCHEM\2\DATA\052506-F\F56124.D\ECD1A.CH Vial: 44
Signal #2 : C:\HPCHEM\2\DATA\052506-F\F56124.D\ECD2B.CH
Acq On : 26 May 2006 9:45 am Operator:
Sample : 56617-20, A/C Inst : Instrumen
Misc : SOIL Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 26 11:18 2006 Quant Results File: PBQ05166.RES

Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 10:06:32 2006
Response via : Multiple Level Calibration
DataAcq Meth : PSTQUICK.M

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film





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Portland ME 04102

May 26, 2006

CLIENT SAMPLE ID

Project Name: GP

Project Number: 218736

Field Sample ID: SS-10

SAMPLE DATA

Lab Sample ID: 56617-21
Matrix: Solid
Percent Solid: 79
Dilution Factor: 1.2
Collection Date: 05/24/06
Lab Receipt Date: 05/25/06
Extraction Date: 05/26/06
Analysis Date: 05/26/06

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit µg/kg	Results µg/kg
PCB-1016	18	U
PCB-1221	18	U
PCB-1232	18	U
PCB-1242	18	U
PCB-1248	18	U
PCB-1254	18	U
PCB-1260	18	**51**

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	85	%
Decachlorobiphenyl	82	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082.

Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3545.

COMMENTS: Results are expressed on a dry weight basis.

PCB Report

Authorized signature Melissa Hall

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PCB CONFIRMATION RESULTS

SDG: 56617

Project No: Site ID: Location: ME

Instrument ID: F GC Columns: RTX-CLPesticidesI ID: 0.32(mm)
RTX-CLPesticidesII

Sample ID: 56617-21

COMPOUND	RESPONSE #1 ug/kg	RESPONSE #2 ug/kg	DIFFERENCE %	#	% DIFFERENCE Acceptance Limit
PCB 1260	51	56	9		40

Column to be used to flag % Difference with an asterisk

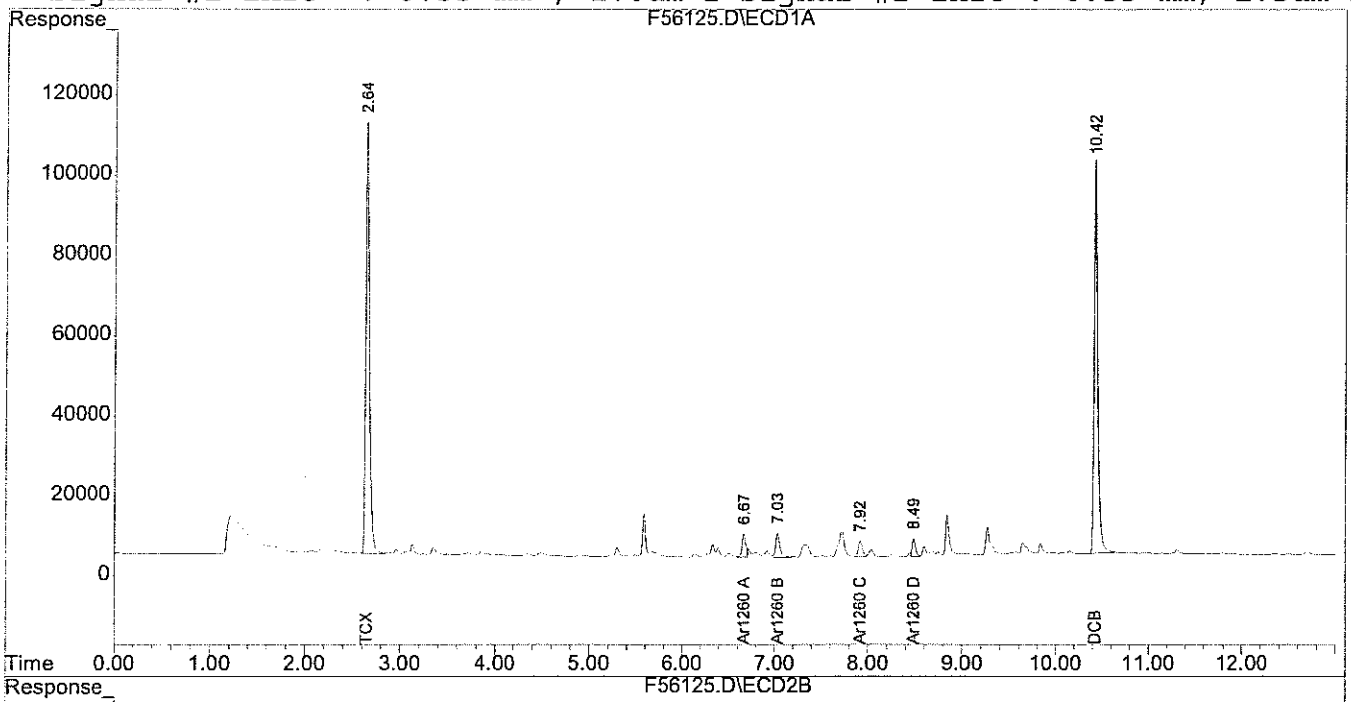
* Values outside of QC limits

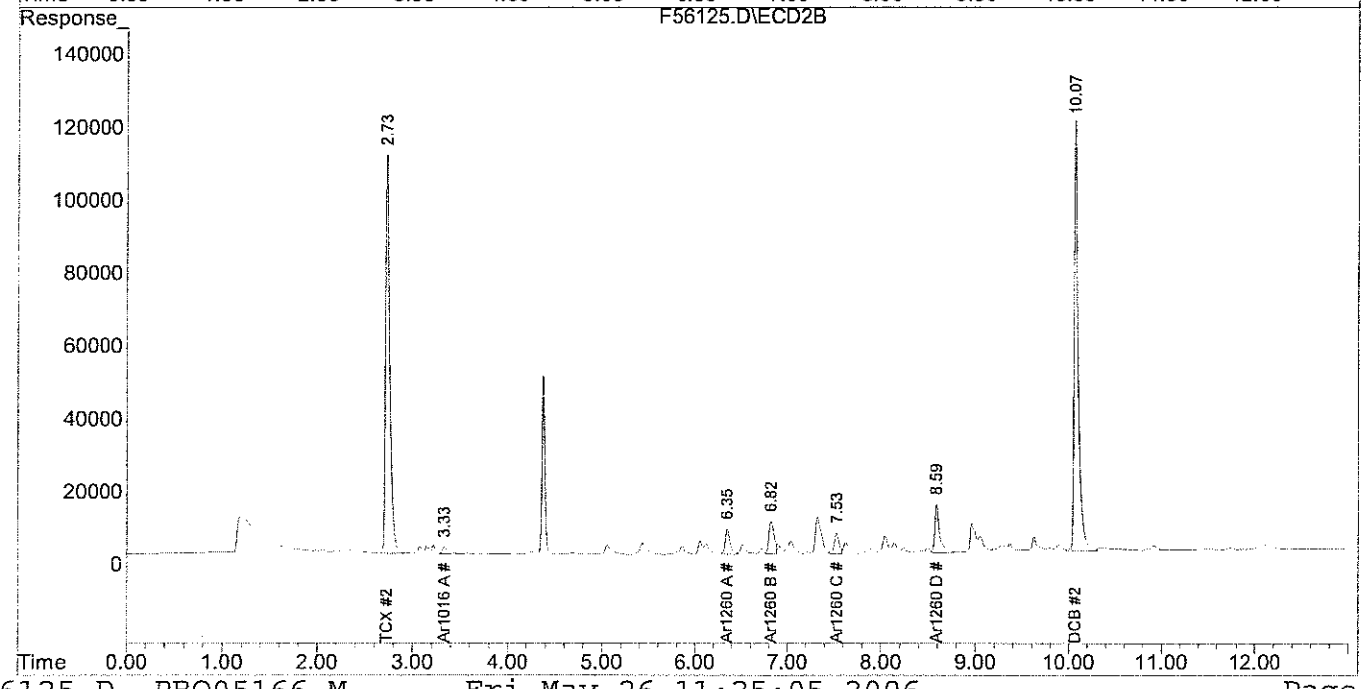
Quantitation Report (QT Reviewed)

Signal #1 : C:\HPCHEM\2\DATA\052506-F\F56125.D\ECD1A.CH Vial: 45
Signal #2 : C:\HPCHEM\2\DATA\052506-F\F56125.D\ECD2B.CH
Acq On : 26 May 2006 10:00 am Operator:
Sample : 56617-21, A/C Inst : Instrumen
Misc : SOIL Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 26 11:24 2006 Quant Results File: PBQ05166.RES

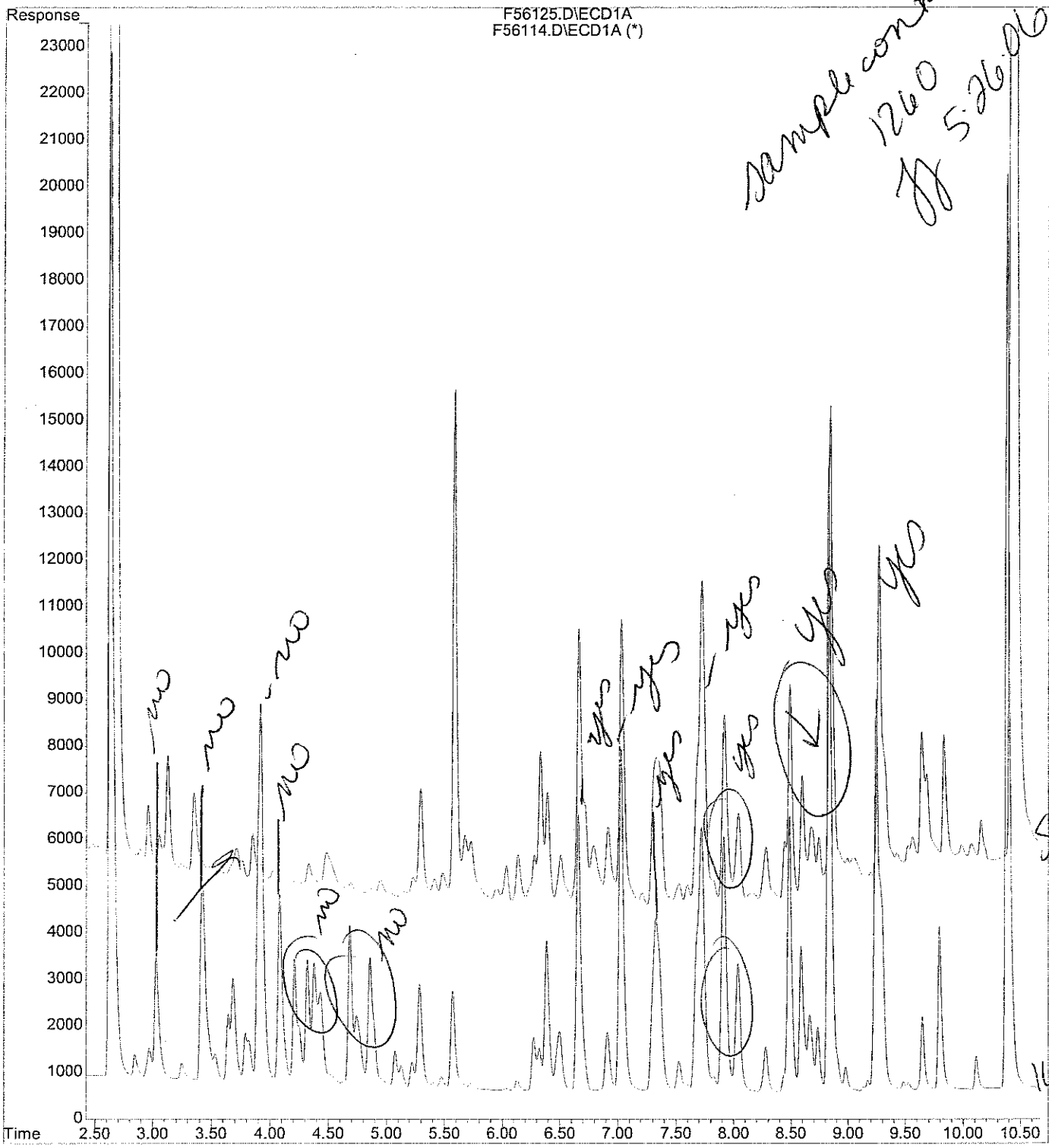
Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 10:06:32 2006
Response via : Multiple Level Calibration
DataAcq Meth : PSTQUICK.M

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film





File : C:\HPCHEM\2\DATA\052506-F\F56125.D
Operator :
Acquired : 26 May 2006 10:00 am using AcqMethod PSTQUICK.M
Instrument : Instrumen
Sample Name: 56617-21, A/C
Misc Info : SOIL
Vial Number: 45



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Portland ME 04102

May 26, 2006

CLIENT SAMPLE ID

Project Name: GP

Project Number: 218736

Field Sample ID: SS-11

SAMPLE DATA

Lab Sample ID:	56617-22
Matrix:	Solid
Percent Solid:	88
Dilution Factor:	1.1
Collection Date:	05/24/06
Lab Receipt Date:	05/25/06
Extraction Date:	05/26/06
Analysis Date:	05/26/06

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit µg/kg	Results µg/kg
PCB-1016	17	U
PCB-1221	17	U
PCB-1232	17	U
PCB-1242	17	U
PCB-1248	17	U
PCB-1254	17	U
PCB-1260	17	65

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	92	%
Decachlorobiphenyl	89	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082.

Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3545.

COMMENTS: Results are expressed on a dry weight basis.

PCB Report

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PCB CONFIRMATION RESULTS

SDG: 56617

Project No: Site ID: Location: ME

Instrument ID: F GC Columns: RTX-CLPesticidesI ID: 0.32(mm)
RTX-CLPesticidesII

Sample ID: 56617-22

COMPOUND	RESPONSE #1 ug/kg	RESPONSE #2 ug/kg	DIFFERENCE %	#	% DIFFERENCE Acceptance Limit
PCB 1260	65	72	11		40

Column to be used to flag % Difference with an asterisk

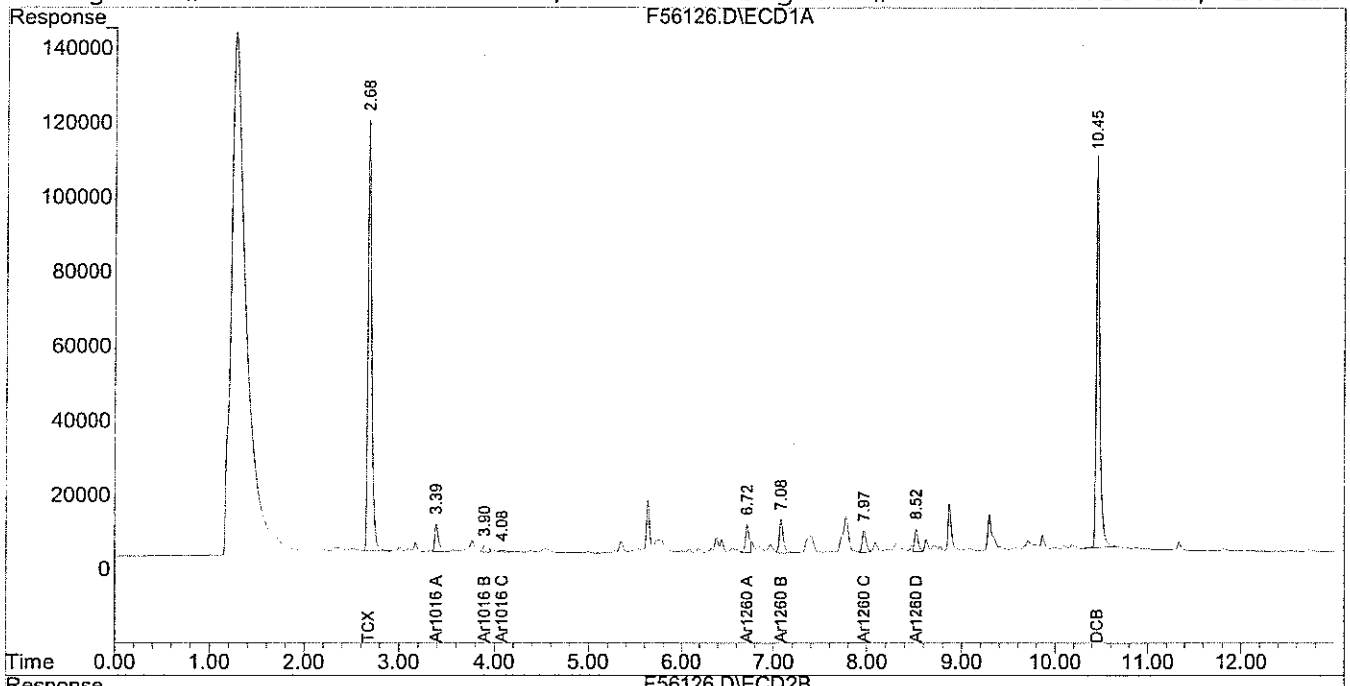
* Values outside of QC limits

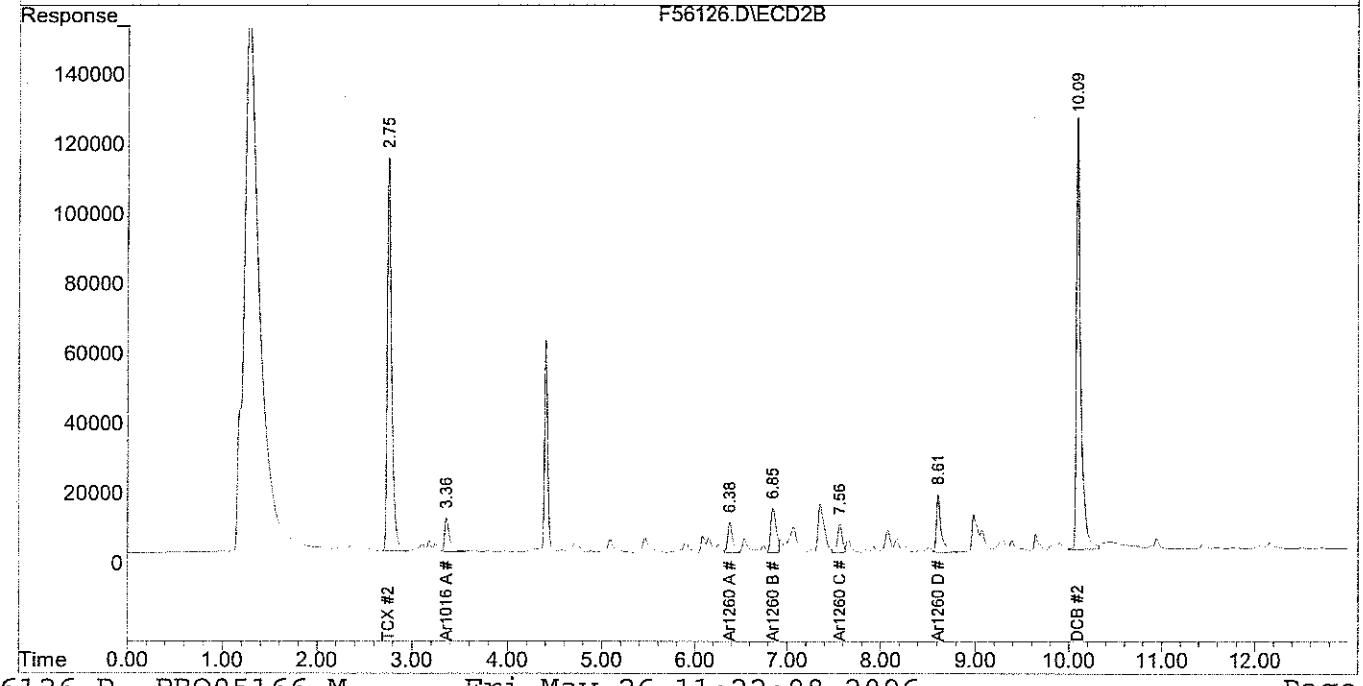
Quantitation Report (QT Reviewed)

Signal #1 : C:\HPCHEM\2\DATA\052506-F\F56126.D\ECD1A.CH Vial: 46
Signal #2 : C:\HPCHEM\2\DATA\052506-F\F56126.D\ECD2B.CH
Acq On : 26 May 2006 10:27 am Operator:
Sample : 56617-22, A/C Inst : Instrumen
Misc : SOIL Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 26 11:21 2006 Quant Results File: PBQ05166.RES

Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 10:06:32 2006
Response via : Multiple Level Calibration
DataAcq Meth : PSTQUICK.M

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film





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File        : C:\HPCHEM\2\DATA\052506-F\F56126.D
Operator    :
Acquired    : 26 May 2006  10:27 am using AcqMethod PSTQUICK.M
Instrument  :   Instrumen
Sample Name: 56617-22, A/C
Misc Info   : SOIL
Vial Number: 46
```

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May 26, 2006

CLIENT SAMPLE ID

Project Name: GP

Project Number: 218736

Field Sample ID: SS-12

SAMPLE DATA

Lab Sample ID: 56617-23
Matrix: Solid
Percent Solid: 90
Dilution Factor: 1.1
Collection Date: 05/24/06
Lab Receipt Date: 05/25/06
Extraction Date: 05/26/06
Analysis Date: 05/26/06

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit µg/kg	Results µg/kg
PCB-1016	17	U
PCB-1221	17	U
PCB-1232	17	U
PCB-1242	17	U
PCB-1248	17	U
PCB-1254	17	U
PCB-1260	17	**49**

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	90	%
Decachlorobiphenyl	84	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082.

Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3545.

COMMENTS: Results are expressed on a dry weight basis.

PCB Report

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PCB CONFIRMATION RESULTS

SDG: 56617

Project No: Site ID: Location: ME

Instrument ID: F GC Columns: RTX-CLPesticidesI ID: 0.32(mm)
RTX-CLPesticidesII

Sample ID: 56617-23

COMPOUND	RESPONSE #1 ug/kg	RESPONSE #2 ug/kg	DIFFERENCE %	#	% DIFFERENCE Acceptance Limit
PCB 1260	49	50	3		40

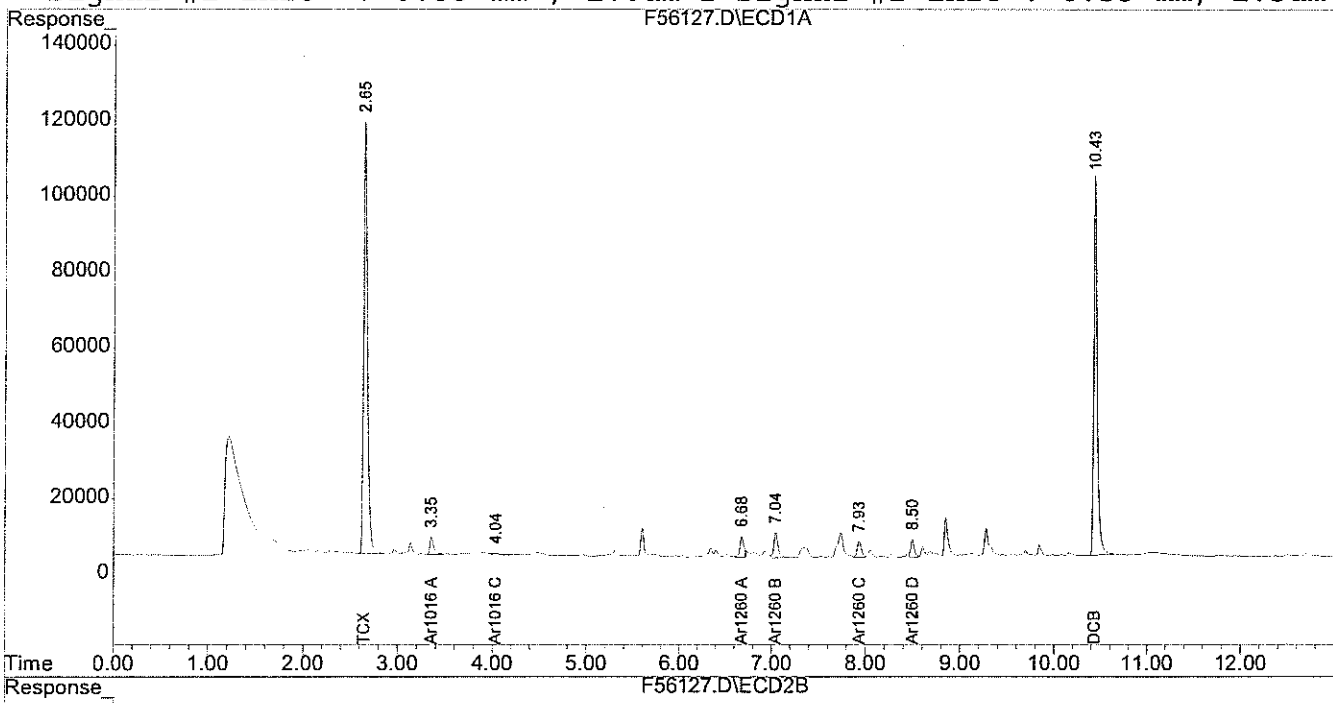
Column to be used to flag % Difference with an asterisk

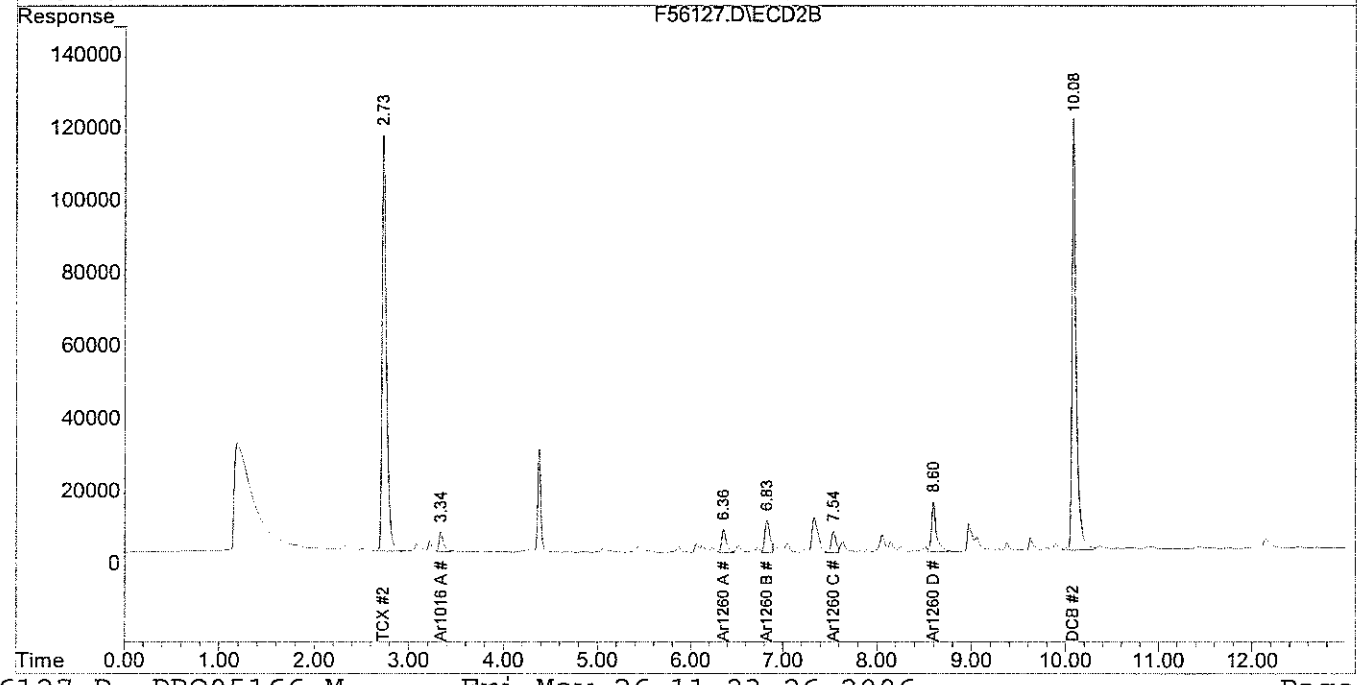
* Values outside of QC limits

Signal #1 : C:\HPCHEM\2\DATA\052506-F\F56127.D\ECD1A.CH Vial: 47
Signal #2 : C:\HPCHEM\2\DATA\052506-F\F56127.D\ECD2B.CH
Acq On : 26 May 2006 10:42 am Operator:
Sample : 56617-23, A/C Inst : Instrumen
Misc : SOIL Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 26 11:23 2006 Quant Results File: PBQ05166.RES

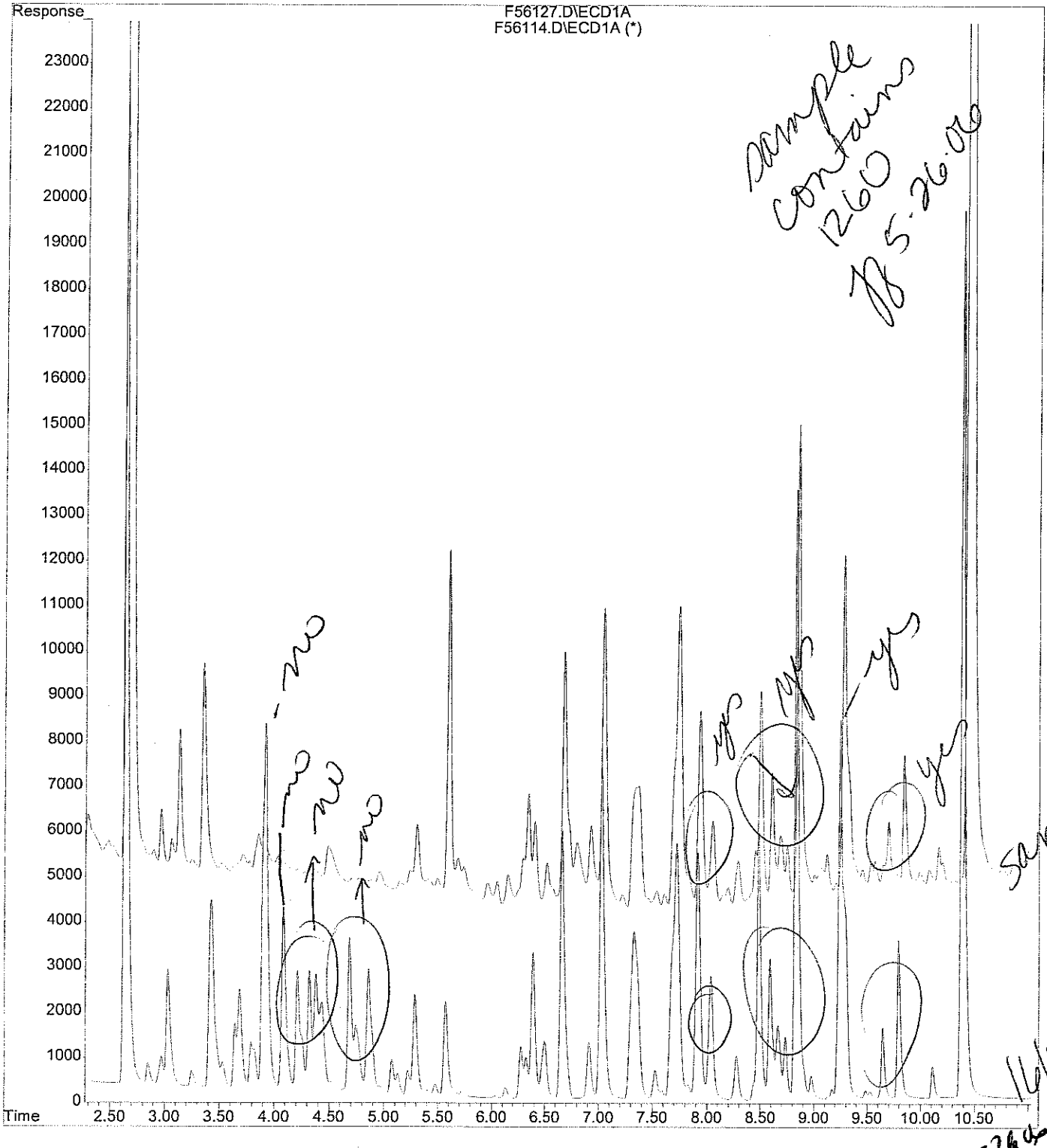
Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 10:06:32 2006
Response via : Multiple Level Calibration
DataAcq Meth : PSTQUICK.M

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film





File : C:\HPCHEM\2\DATA\052506-F\F56127.D
Operator :
Acquired : 26 May 2006 10:42 am using AcqMethod PSTQUICK.M
Instrument : Instrumen
Sample Name: 56617-23, A/C
Misc Info : SOIL
Vial Number: 47



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May 26, 2006

CLIENT SAMPLE ID

Project Name: GP

Project Number: 218736

Field Sample ID: W-1

SAMPLE DATA

Lab Sample ID: 56617-24
Matrix: Wipe
Percent Solid: N/A
Dilution Factor: 1.0
Collection Date: 05/24/06
Lab Receipt Date: 05/25/06
Extraction Date: 05/25/06
Analysis Date: 05/25/06

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit ug/wipe	Results ug/wipe
PCB-1016	0.5	U
PCB-1221	0.5	U
PCB-1232	0.5	U
PCB-1242	0.5	U
PCB-1248	0.5	U
PCB-1254	0.5	U
PCB-1260	0.5	1.5

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	93	%
Decachlorobiphenyl	103	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8081/82.

COMMENTS:

Authorized signature

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PCB CONFIRMATION RESULTS

SDG: 56617

Project No: Site ID: Location: ME

Instrument ID: F GC Columns: RTX-CLPesticidesI ID: 0.32(mm)
RTX-CLPesticidesII

Sample ID: 56617-24

COMPOUND	RESPONSE #1 ug/wipe	RESPONSE #2 ug/wipe	DIFFERENCE %	#	% DIFFERENCE Acceptance Limit
PCB 1260	1.5	1.8	23		40

Column to be used to flag % Difference with an asterisk

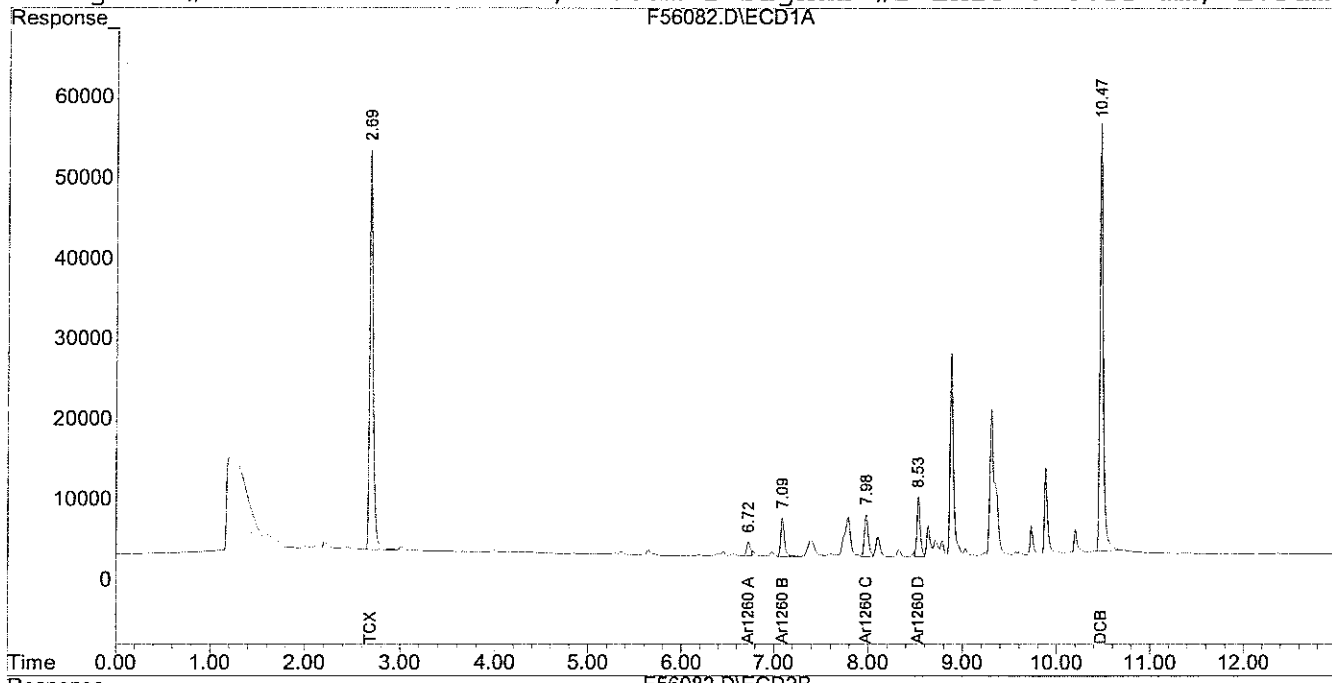
* Values outside of QC limits

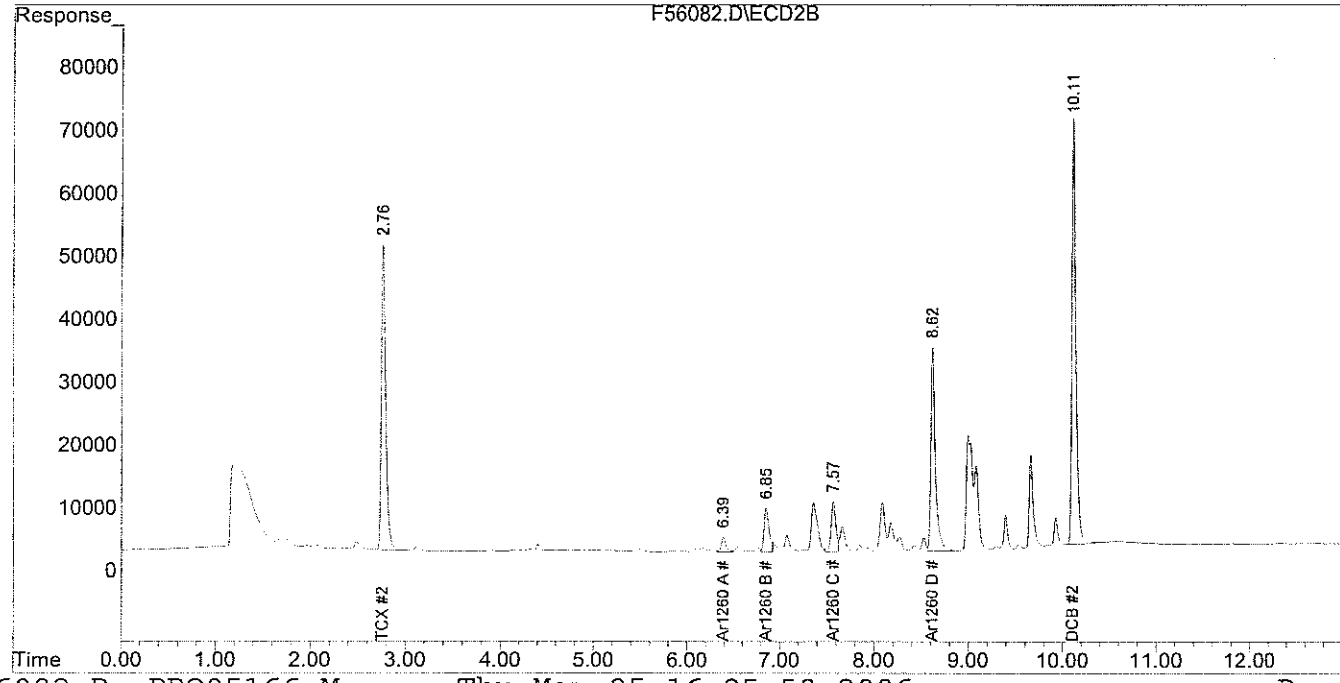
Quantitation Report (QT Reviewed)

Signal #1 : C:\HPCHEM\2\DATA\052506-F\F56082.D\ECD1A.CH Vial: 5
Signal #2 : C:\HPCHEM\2\DATA\052506-F\F56082.D\ECD2B.CH
Acq On : 25 May 2006 4:08 pm Operator:
Sample : 56617-24, A/C Inst : Instrumen
Misc : SOIL Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 25 16:25 2006 Quant Results File: PBQ05166.RES

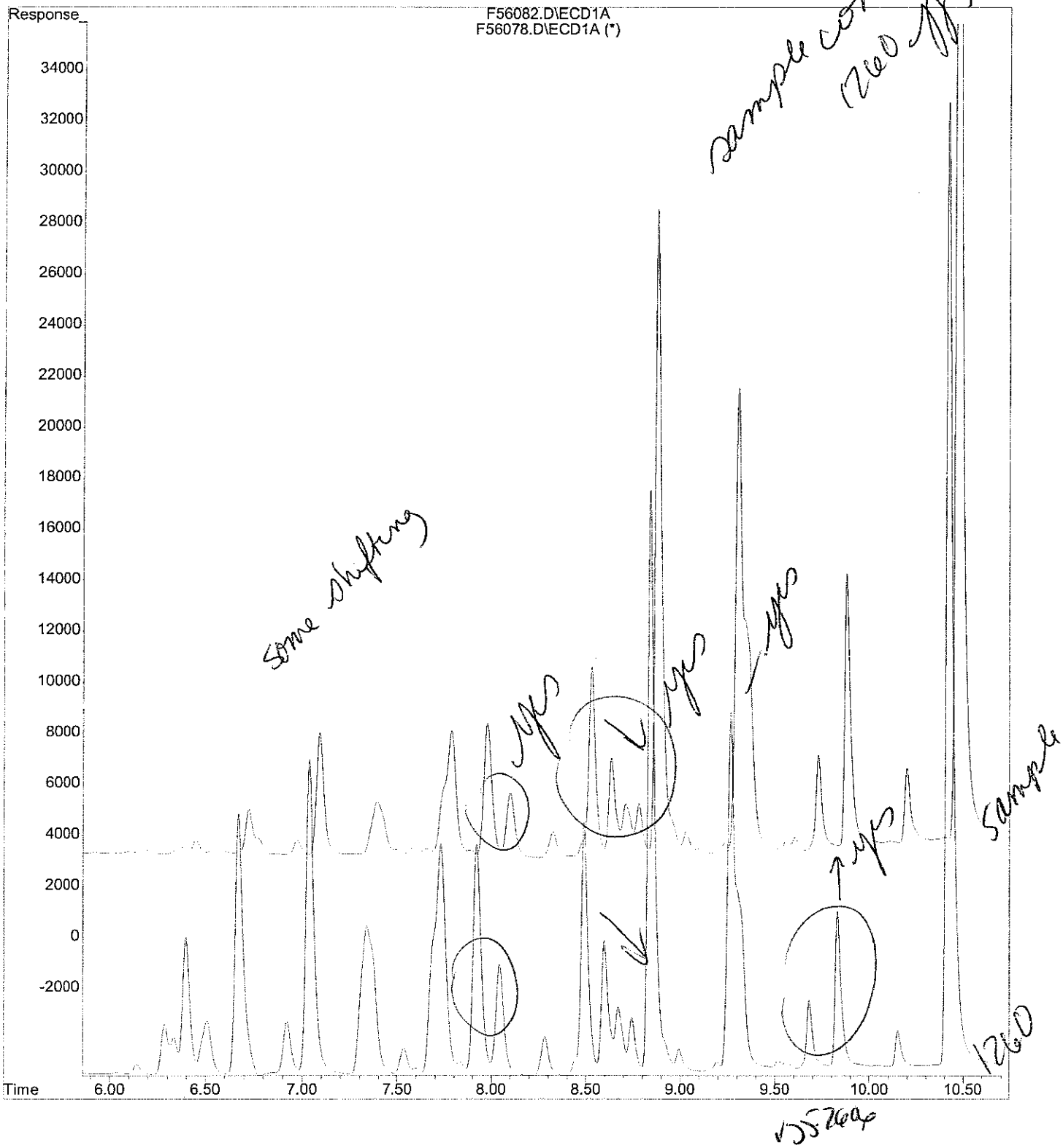
Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 10:06:32 2006
Response via : Multiple Level Calibration
DataAcq Meth : PSTQUICK.M

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film





File : C:\HPCHEM\2\DATA\052506-F\F56082.D
Operator :
Acquired : 25 May 2006 4:08 pm using AcqMethod PSTQUICK.M
Instrument : Instrumen
Sample Name: 56617-24, A/C
Misc Info : SOIL
Vial Number: 5



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May 26, 2006

CLIENT SAMPLE ID

Project Name: GP

Project Number: 218736

Field Sample ID: W-2

SAMPLE DATA

Lab Sample ID: 56617-25
Matrix: Wipe
Percent Solid: N/A
Dilution Factor: 1.0
Collection Date: 05/24/06
Lab Receipt Date: 05/25/06
Extraction Date: 05/25/06
Analysis Date: 05/25/06

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit ug/wipe	Results ug/wipe
PCB-1016	0.5	U
PCB-1221	0.5	U
PCB-1232	0.5	U
PCB-1242	0.5	U
PCB-1248	0.5	U
PCB-1254	0.5	U
PCB-1260	0.5	**4.2**

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	98	%
Decachlorobiphenyl	108	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8081/82.

COMMENTS:

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PCB CONFIRMATION RESULTS

SDG: 56617

Project No: Site ID: Location: ME

Instrument ID: F GC Columns: RTX-CLPesticidesI ID: 0.32(mm)
RTX-CLPesticidesII

Sample ID: 56617-25

COMPOUND	RESPONSE #1 ug/wipe	RESPONSE #2 ug/wipe	DIFFERENCE %	#	% DIFFERENCE Acceptance Limit
PCB 1260	4.2	6.0	34		40

Column to be used to flag % Difference with an asterisk

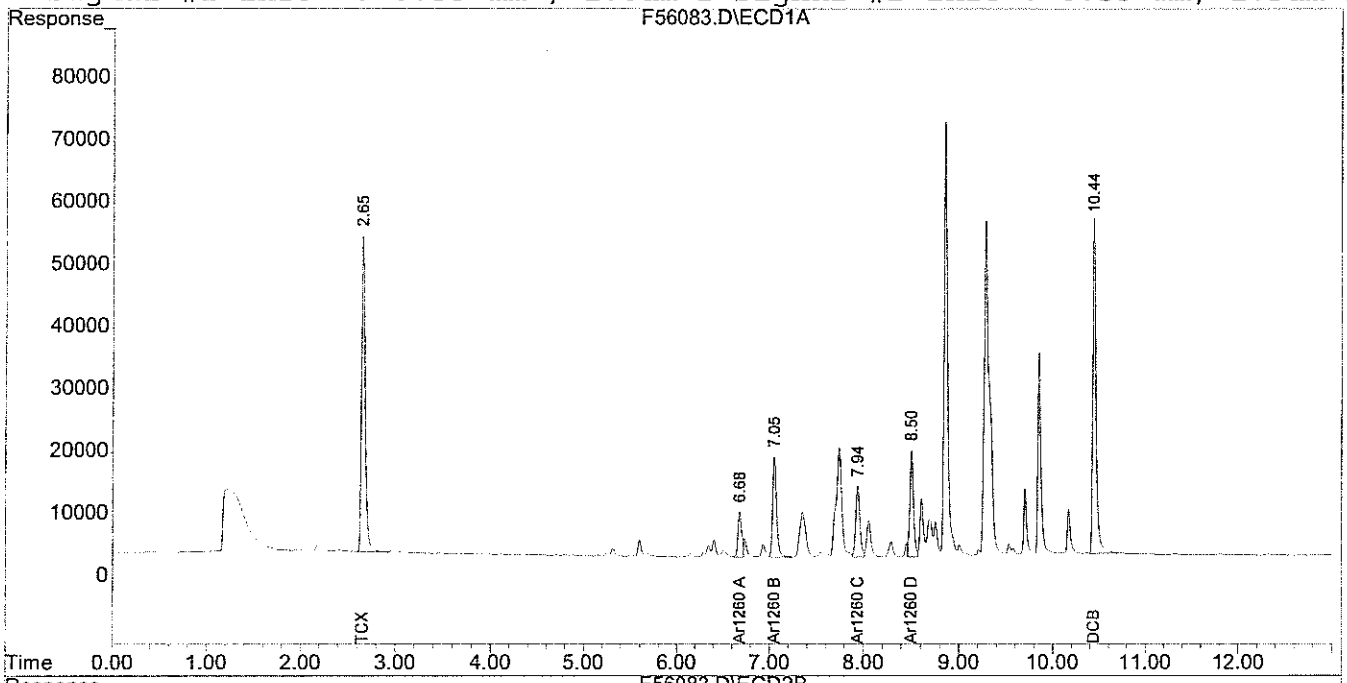
* Values outside of QC limits

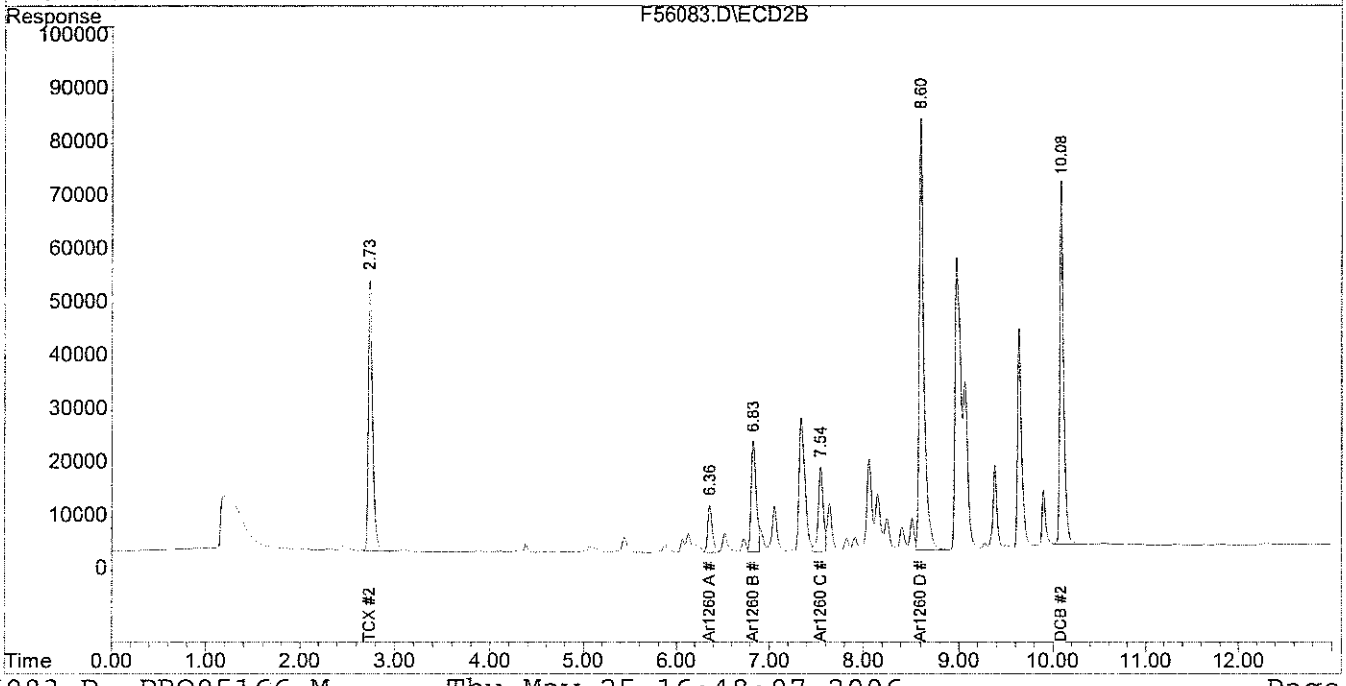
Quantitation Report (QT Reviewed)

Signal #1 : C:\HPCHEM\2\DATA\052506-F\F56083.D\ECD1A.CH Vial: 6
Signal #2 : C:\HPCHEM\2\DATA\052506-F\F56083.D\ECD2B.CH
Acq On : 25 May 2006 4:23 pm Operator:
Sample : 56617-25, A/C Inst : Instrumen
Misc : SOIL Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 25 16:48 2006 Quant Results File: PBQ05166.RES

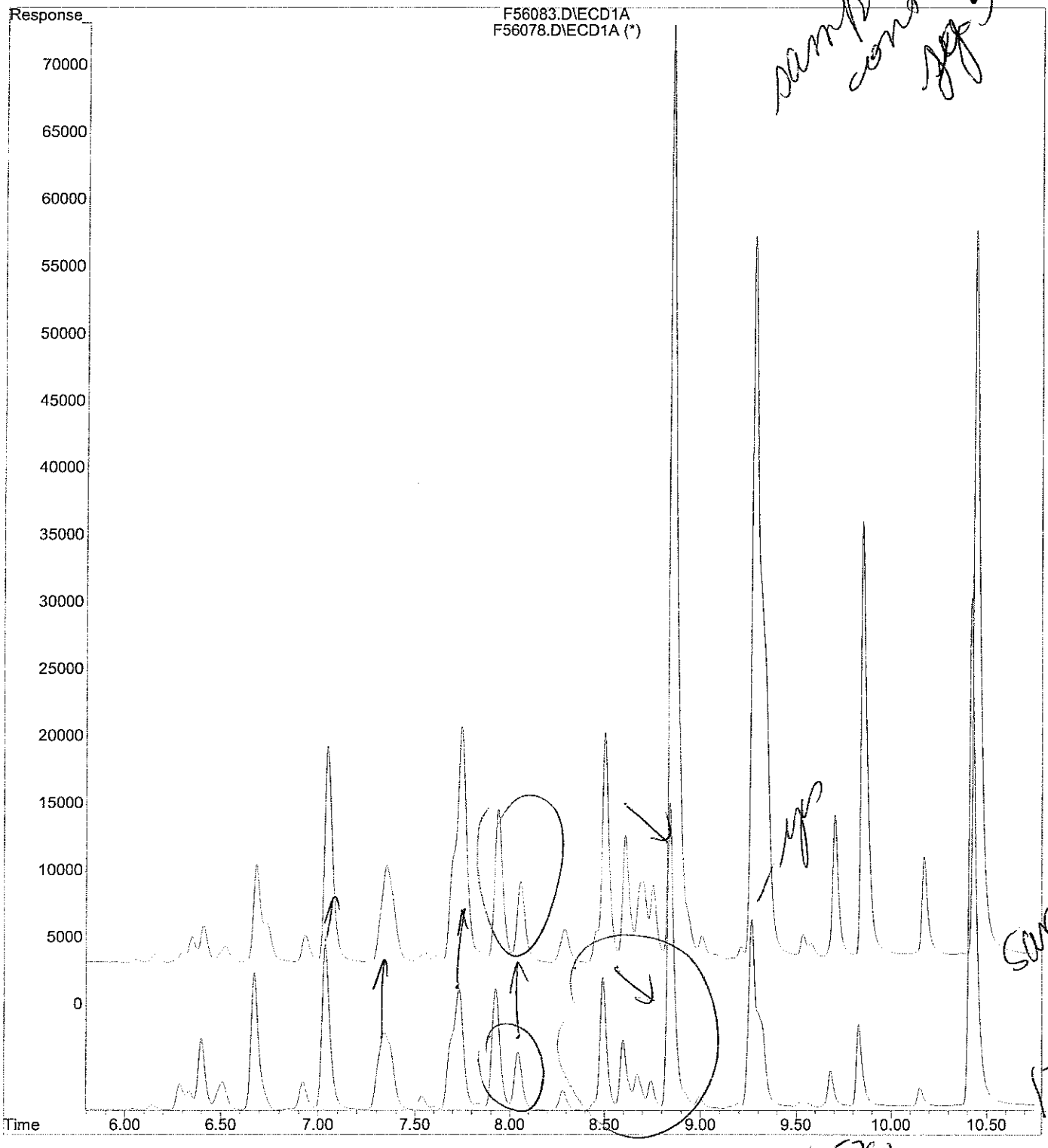
Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 10:06:32 2006
Response via : Multiple Level Calibration
DataAcq Meth : PSTQUICK.M

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film





File : C:\HPCHEM\2\DATA\052506-F\F56083.D
Operator :
Acquired : 25 May 2006 4:23 pm using AcqMethod PSTQUICK.M
Instrument : Instrumen
Sample Name: 56617-25, A/C
Misc Info : SOIL
Vial Number: 6



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May 26, 2006

CLIENT SAMPLE ID

Project Name: GP

Project Number: 218736

Field Sample ID: W-3

SAMPLE DATA

Lab Sample ID: 56617-26
Matrix: Wipe
Percent Solid: N/A
Dilution Factor: 1.0
Collection Date: 05/24/06
Lab Receipt Date: 05/25/06
Extraction Date: 05/25/06
Analysis Date: 05/25/06

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit ug/wipe	Results ug/wipe
PCB-1016	0.5	U
PCB-1221	0.5	U
PCB-1232	0.5	U
PCB-1242	0.5	U
PCB-1248	0.5	U
PCB-1254	0.5	U
PCB-1260	0.5	**6.1**

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	98	%
Decachlorobiphenyl	108	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8081/82.

COMMENTS:

PCB Report

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PCB CONFIRMATION RESULTS

SDG: 56617

Project No: Site ID: Location: ME

Instrument ID: F GC Columns: RTX-CLPesticidesI ID: 0.32(mm)
RTX-CLPesticidesII

Sample ID: 56617-26

COMPOUND	RESPONSE #1 ug/wipe	RESPONSE #2 ug/wipe	DIFFERENCE %	#	% DIFFERENCE Acceptance Limit
PCB 1260	6.1	8.2	28		40

Column to be used to flag % Difference with an asterisk

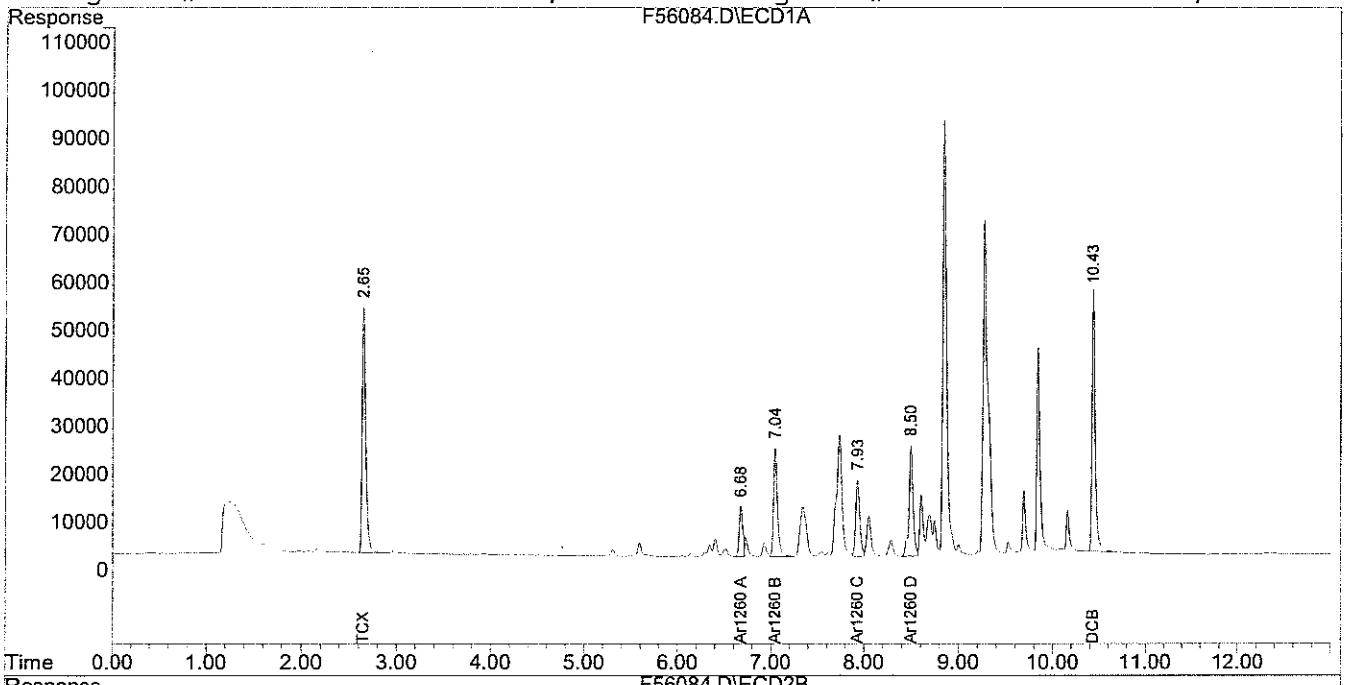
* Values outside of QC limits

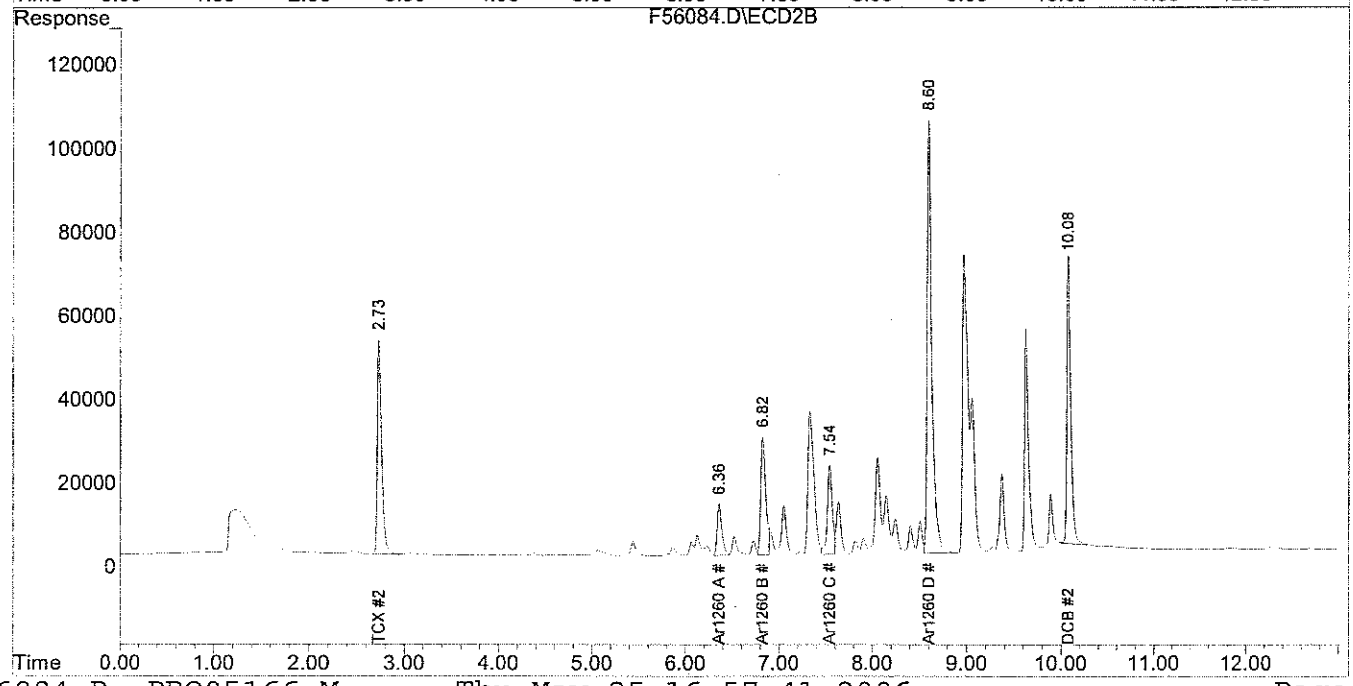
Quantitation Report (QT Reviewed)

Signal #1 : C:\HPCHEM\2\DATA\052506-F\F56084.D\ECD1A.CH Vial: 7
Signal #2 : C:\HPCHEM\2\DATA\052506-F\F56084.D\ECD2B.CH
Acq On : 25 May 2006 4:52 pm Operator:
Sample : 56617-26, A/C Inst : Instrumen
Misc : SOIL Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 25 16:57 2006 Quant Results File: PBQ05166.RES

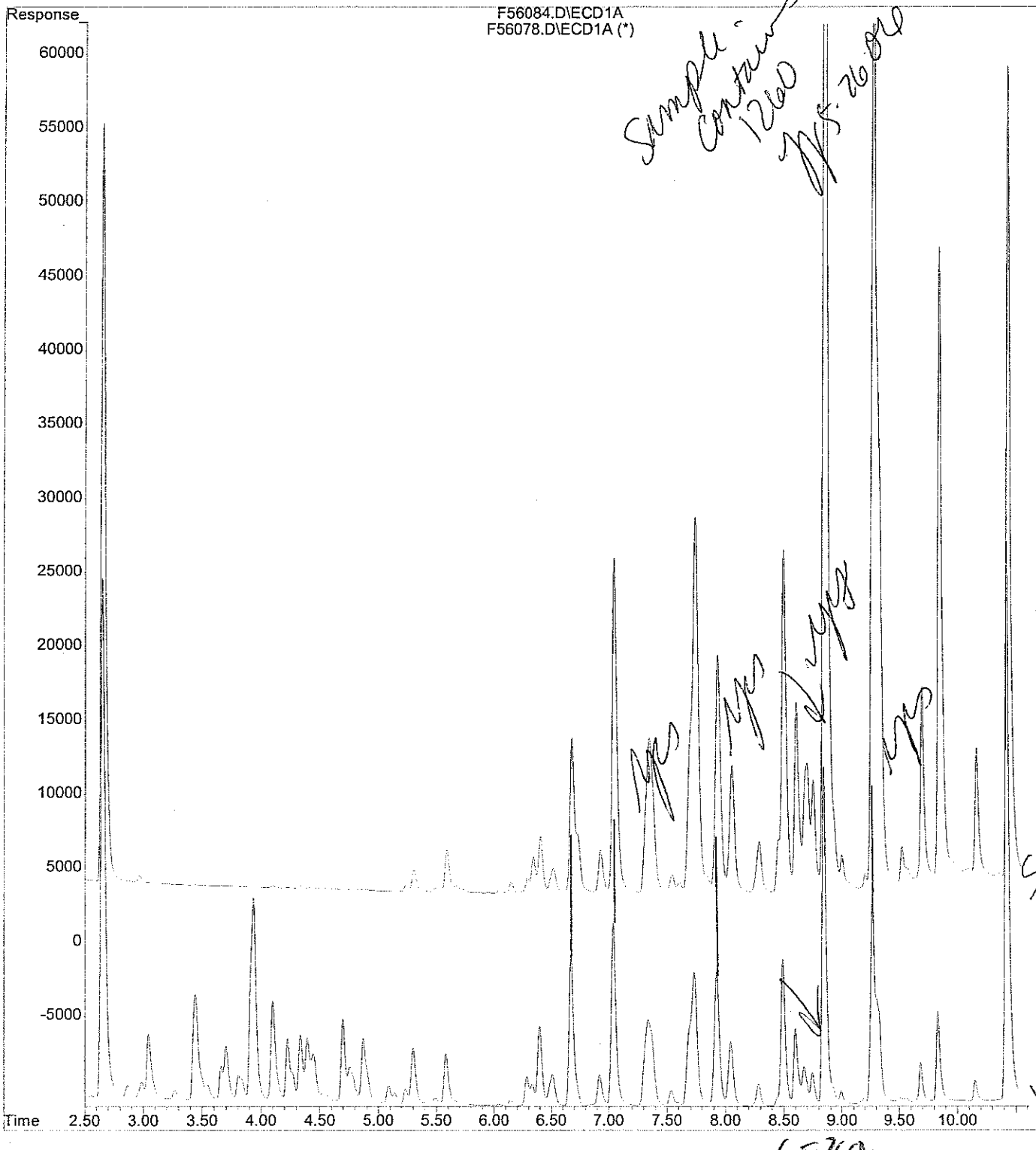
Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 10:06:32 2006
Response via : Multiple Level Calibration
DataAcq Meth : PSTQUICK.M

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film





File : C:\HPCHEM\2\DATA\052506-F\F56084.D
Operator :
Acquired : 25 May 2006 4:52 pm using AcqMethod PSTQUICK.M
Instrument : Instrumen
Sample Name: 56617-26, A/C
Misc Info : SOIL
Vial Number: 7



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May 26, 2006

CLIENT SAMPLE ID

Project Name: GP

Project Number: 218736

Field Sample ID: W-4

SAMPLE DATA

Lab Sample ID:	56617-27
Matrix:	Wipe
Percent Solid:	N/A
Dilution Factor:	1.0
Collection Date:	05/24/06
Lab Receipt Date:	05/25/06
Extraction Date:	05/25/06
Analysis Date:	05/25/06

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit ug/wipe	Results ug/wipe
PCB-1016	0.5	U
PCB-1221	0.5	U
PCB-1232	0.5	U
PCB-1242	0.5	U
PCB-1248	0.5	U
PCB-1254	0.5	U
PCB-1260	0.5	U

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	98	%
Decachlorobiphenyl	105	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8081/82.

COMMENTS:

PCB Report

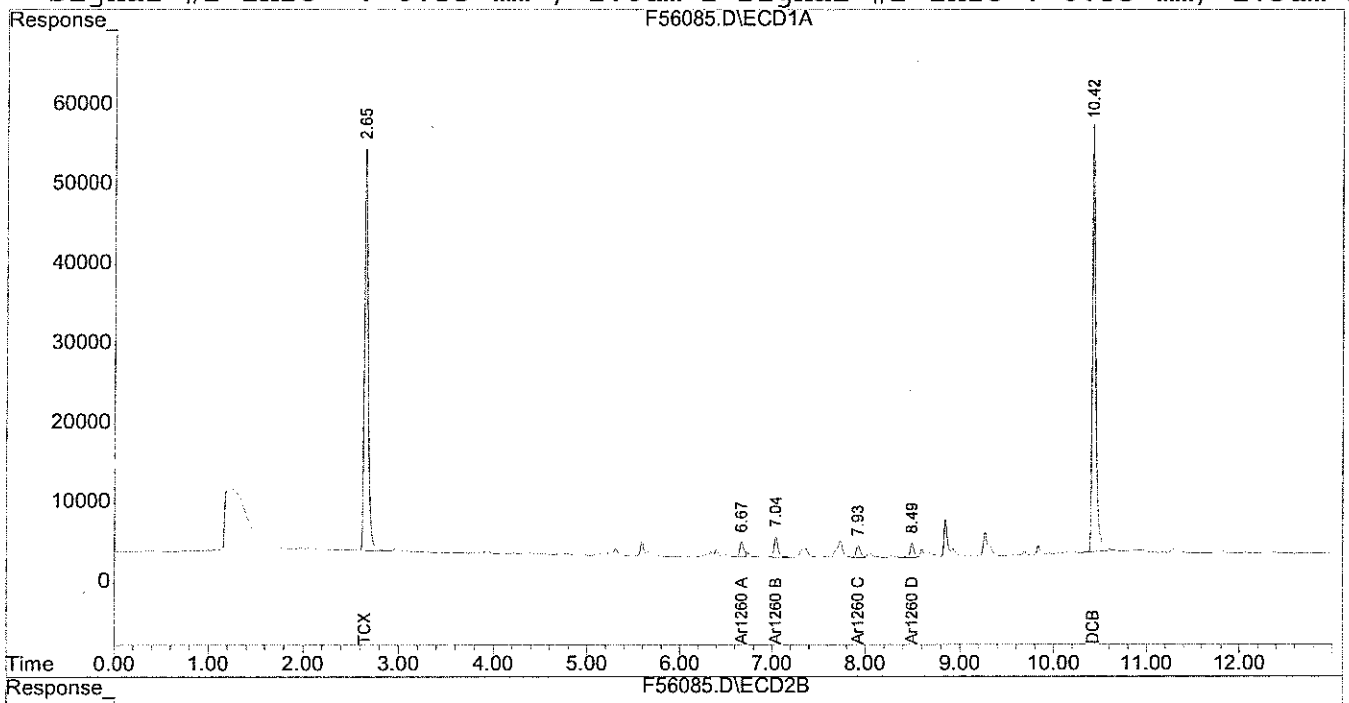
Authorized signature

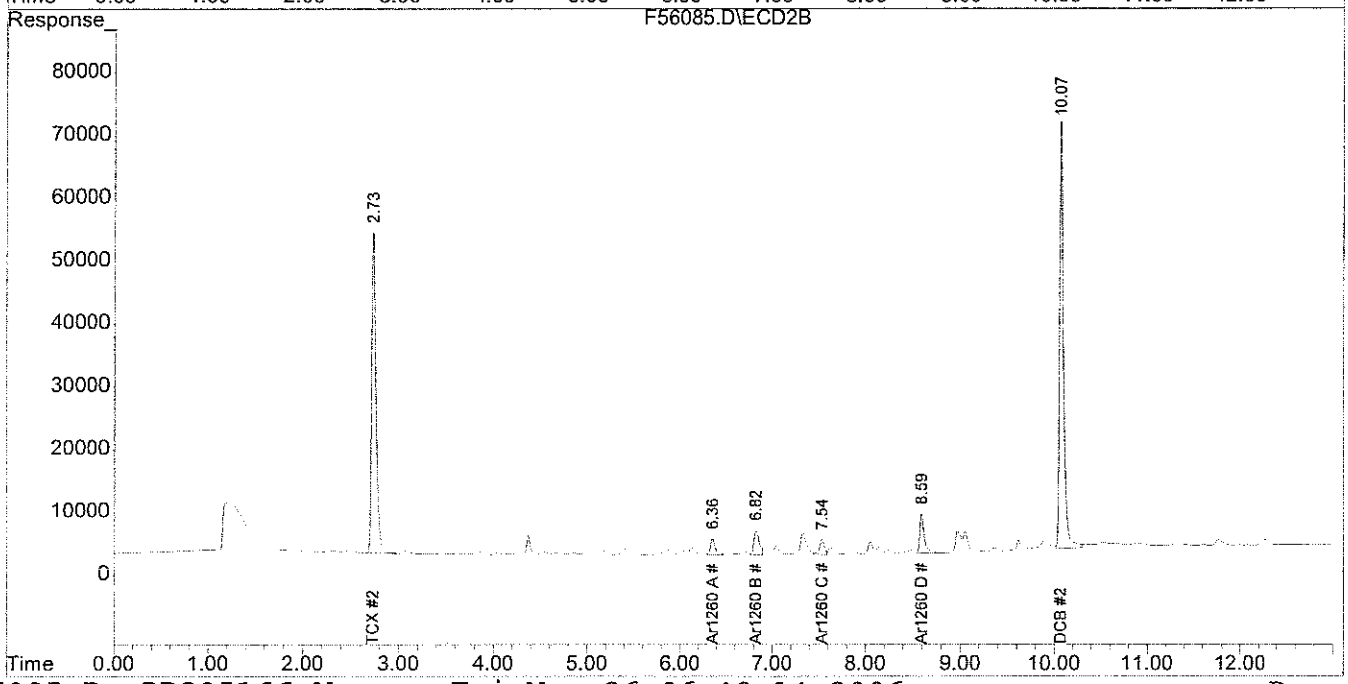
Quantitation Report (QT Reviewed)

Signal #1 : C:\HPCHEM\2\DATA\052506-F\F56085.D\ECD1A.CH Vial: 8
Signal #2 : C:\HPCHEM\2\DATA\052506-F\F56085.D\ECD2B.CH
Acq On : 25 May 2006 5:07 pm Operator:
Sample : 56617-27, A/C Inst : Instrumen
Misc : SOIL Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 26 6:49 2006 Quant Results File: PBQ05166.RES

Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 10:06:32 2006
Response via : Multiple Level Calibration
DataAcq Meth : PSTQUICK.M

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film





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Portland ME 04102

May 30, 2006

CLIENT SAMPLE ID

Project Name: GP

Project Number: 218736

Field Sample ID: W-5

SAMPLE DATA

Lab Sample ID: 56617-28 DL
Matrix: Wipe
Percent Solid: N/A
Dilution Factor: 10
Collection Date: 05/24/06
Lab Receipt Date: 05/25/06
Extraction Date: 05/25/06
Analysis Date: 05/26/06

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit µg/wipe	Results µg/wipe
PCB-1016	5	U
PCB-1221	5	U
PCB-1232	5	U
PCB-1242	5	U
PCB-1248	5	U
PCB-1254	5	U
PCB-1260	5	**107**

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	75	%
Decachlorobiphenyl	50	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8081/82.

Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3550B.

COMMENTS:

PCB Report

Authorized signature Melissa Hall

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PCB CONFIRMATION RESULTS

SDG: 56617

Project No: Site ID: Location: ME

Instrument ID: F GC Columns: RTX-CLPesticidesI ID: 0.32(mm)

RTX-CLPesticidesII

Sample ID: 56617-28

COMPOUND	RESPONSE #1 ug/wipe	RESPONSE #2 ug/wipe	DIFFERENCE %	#	% DIFFERENCE Acceptance Limit
PCB 1260	107	142	28		40

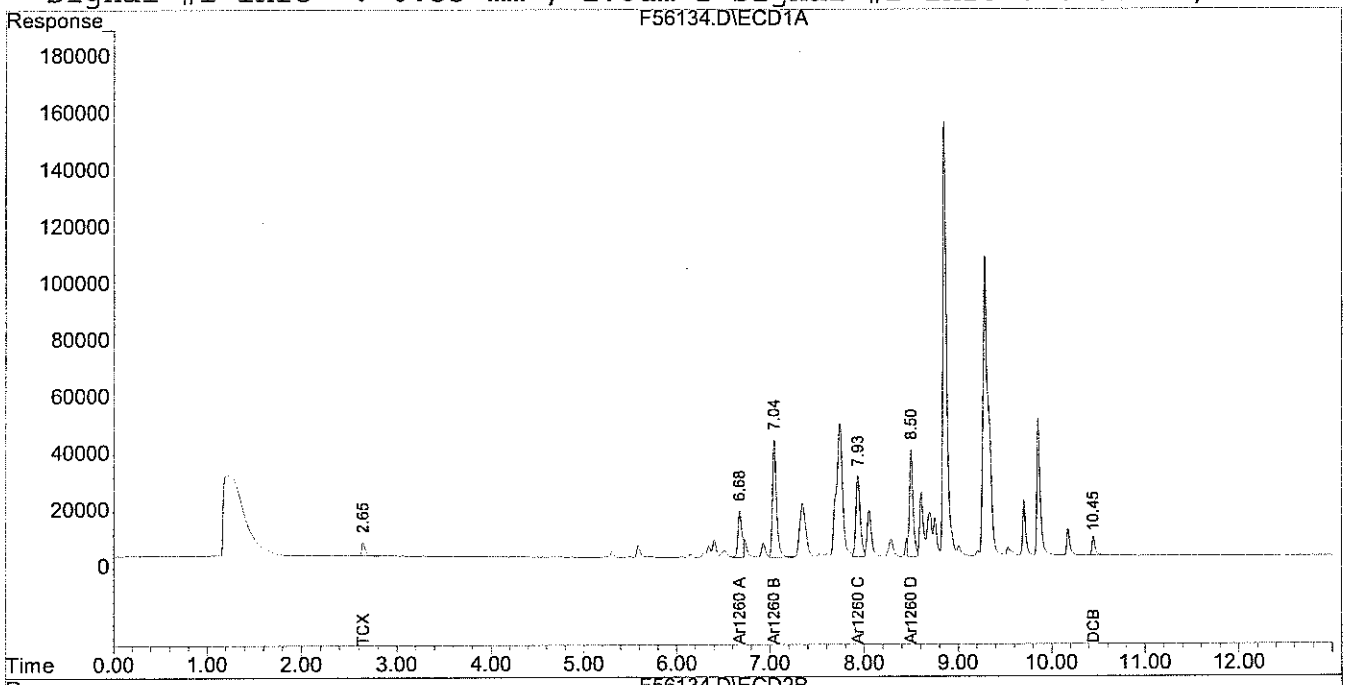
Column to be used to flag % Difference with an asterisk

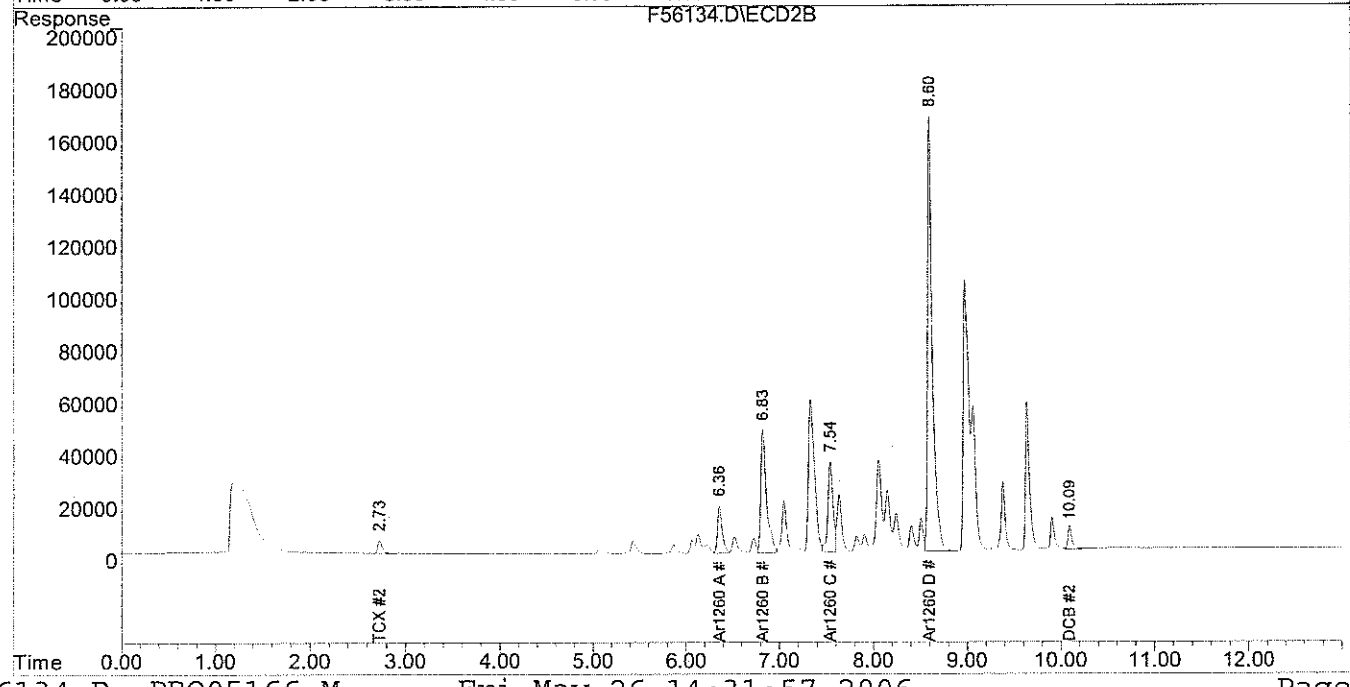
* Values outside of QC limits

Signal #1 : C:\HPCHEM\2\DATA\052606-F\F56134.D\ECD1A.CH Vial: 51
Signal #2 : C:\HPCHEM\2\DATA\052606-F\F56134.D\ECD2B.CH
Acq On : 26 May 2006 2:22 pm Operator:
Sample : 56617-28, 1:10, A/C Inst : Instrumen
Misc : SOIL Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 26 14:31 2006 Quant Results File: PBQ05166.RES

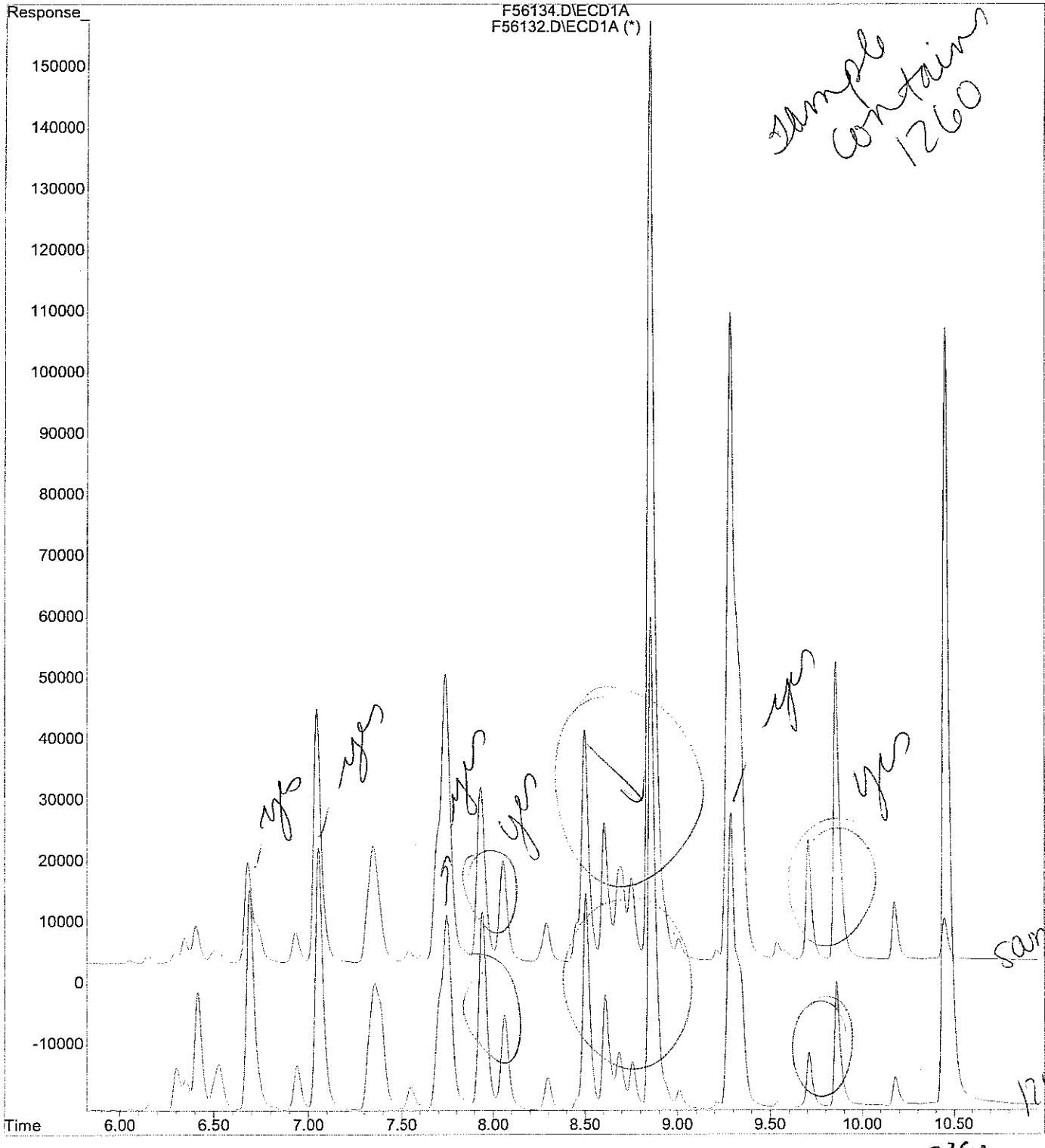
Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 10:06:32 2006
Response via : Multiple Level Calibration
DataAcq Meth : PSTQUICK.M

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film





File : C:\HPCHEM\2\DATA\052606-F\F56134.D
Operator :
Acquired : 26 May 2006 2:22 pm using AcqMethod PSTQUICK.M
Instrument : Instrumen
Sample Name: 56617-28, 1:10, A/C
Misc Info : SOIL
Vial Number: 51



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Mr. Karl Kasper
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41 Hutchins Drive
Portland ME 04102

May 26, 2006

CLIENT SAMPLE ID

Project Name: GP

Project Number: 218736

Field Sample ID: W-6

SAMPLE DATA

Lab Sample ID: 56617-29
Matrix: Wipe
Percent Solid: N/A
Dilution Factor: 1.0
Collection Date: 05/24/06
Lab Receipt Date: 05/25/06
Extraction Date: 05/25/06
Analysis Date: 05/25/06

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit ug/wipe	Results ug/wipe
PCB-1016	0.5	U
PCB-1221	0.5	U
PCB-1232	0.5	U
PCB-1242	0.5	U
PCB-1248	0.5	U
PCB-1254	0.5	U
PCB-1260	0.5	U

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	98	%
Decachlorobiphenyl	110	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8081/82.

COMMENTS:

PCB Report

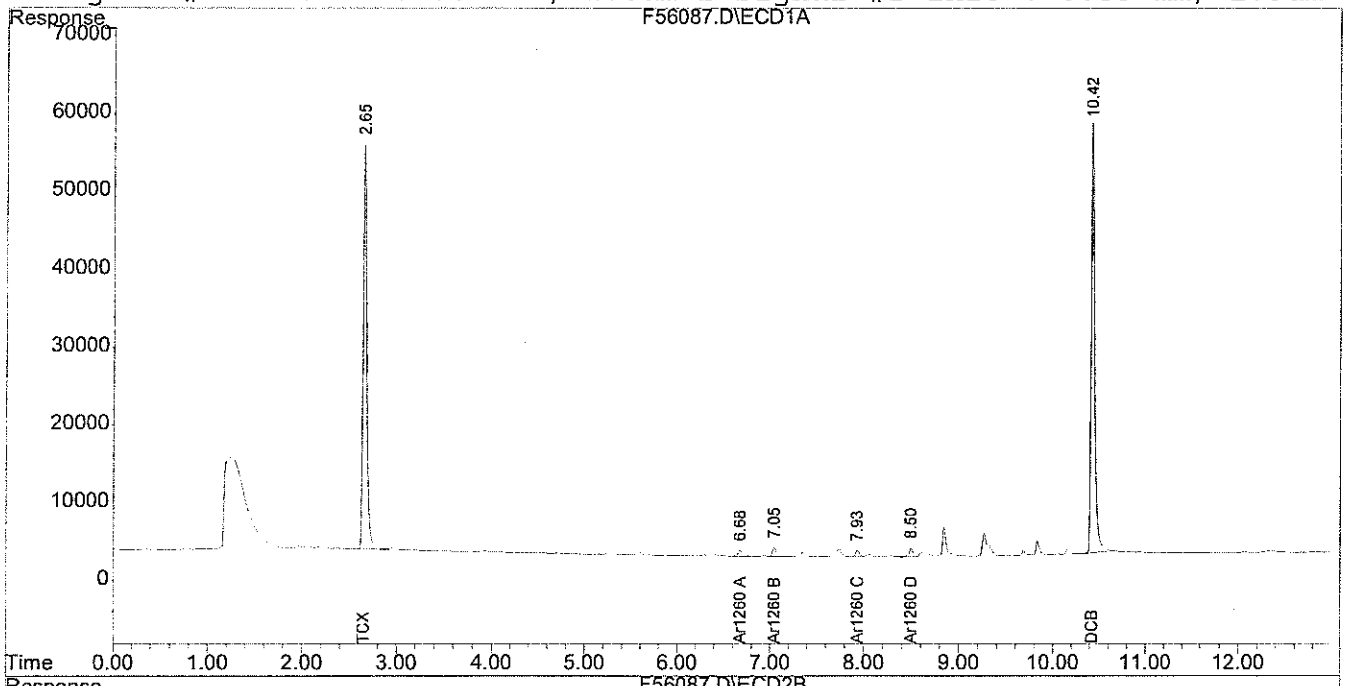
Authorized signature

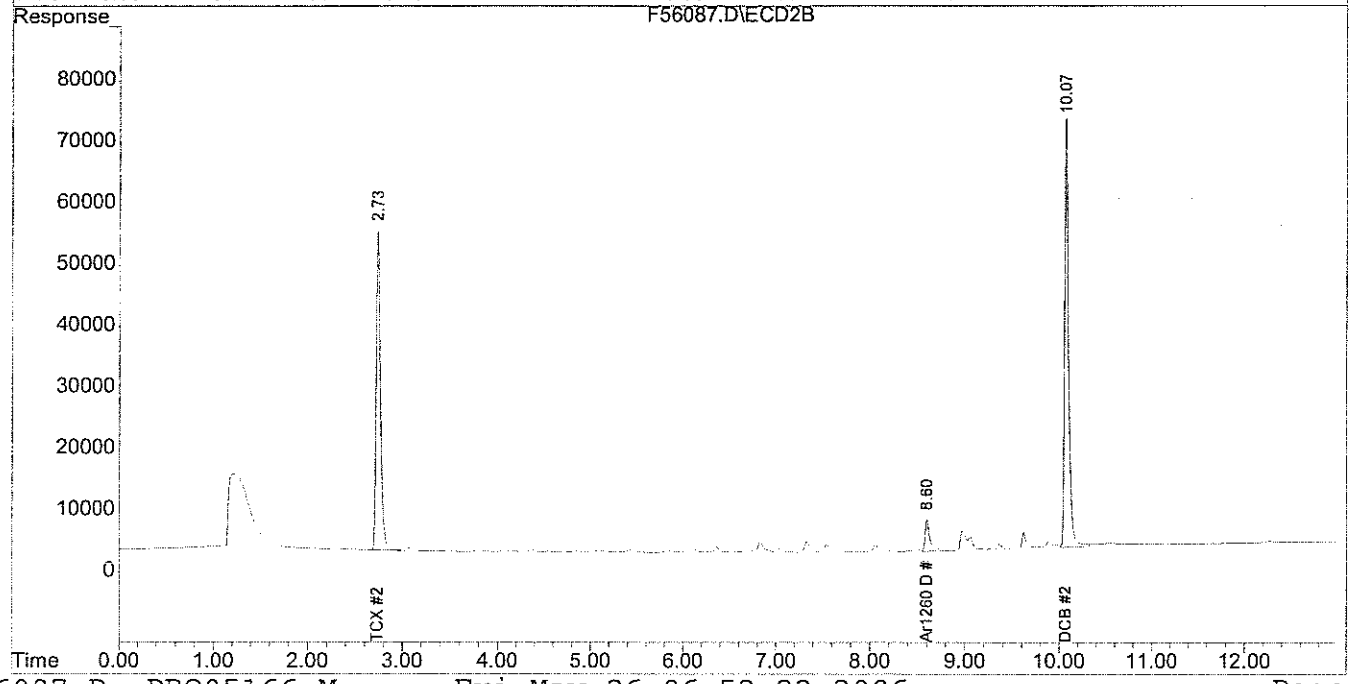
Quantitation Report (QT Reviewed)

Signal #1 : C:\HPCHEM\2\DATA\052506-F\F56087.D\ECD1A.CH Vial: 10
Signal #2 : C:\HPCHEM\2\DATA\052506-F\F56087.D\ECD2B.CH
Acq On : 25 May 2006 5:36 pm Operator:
Sample : 56617-29, A/C Inst : Instrumen
Misc : SOIL Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 26 6:52 2006 Quant Results File: PBQ05166.RES

Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 10:06:32 2006
Response via : Multiple Level Calibration
DataAcq Meth : PSTQUICK.M

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film





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May 26, 2006

CLIENT SAMPLE ID

Project Name: GP

Project Number: 218736

Field Sample ID: W-6 DUP

SAMPLE DATA

Lab Sample ID: 56617-30
Matrix: Wipe
Percent Solid: N/A
Dilution Factor: 1.0
Collection Date: 05/24/06
Lab Receipt Date: 05/25/06
Extraction Date: 05/25/06
Analysis Date: 05/25/06

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit ug/wipe	Results ug/wipe
PCB-1016	0.5	U
PCB-1221	0.5	U
PCB-1232	0.5	U
PCB-1242	0.5	U
PCB-1248	0.5	U
PCB-1254	0.5	U
PCB-1260	0.5	U

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	100	%
Decachlorobiphenyl	108	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8081/82.

COMMENTS:

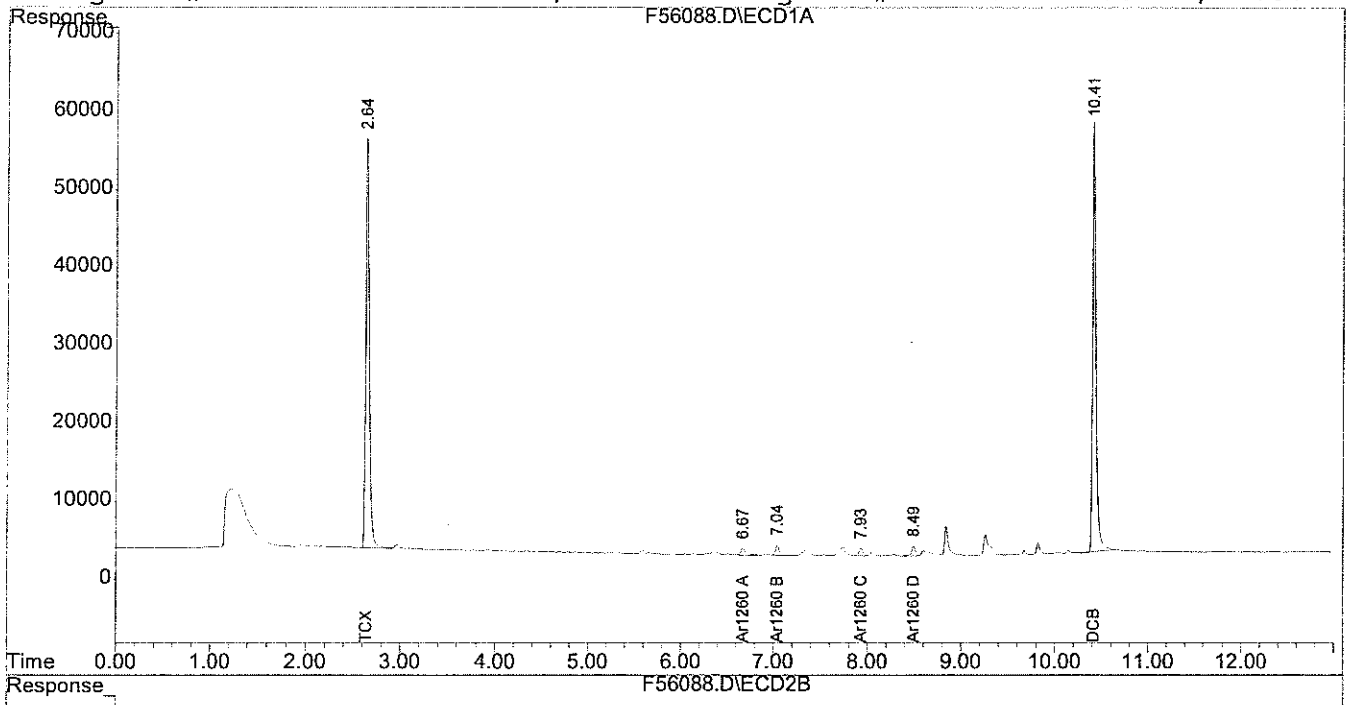
Authorized signature

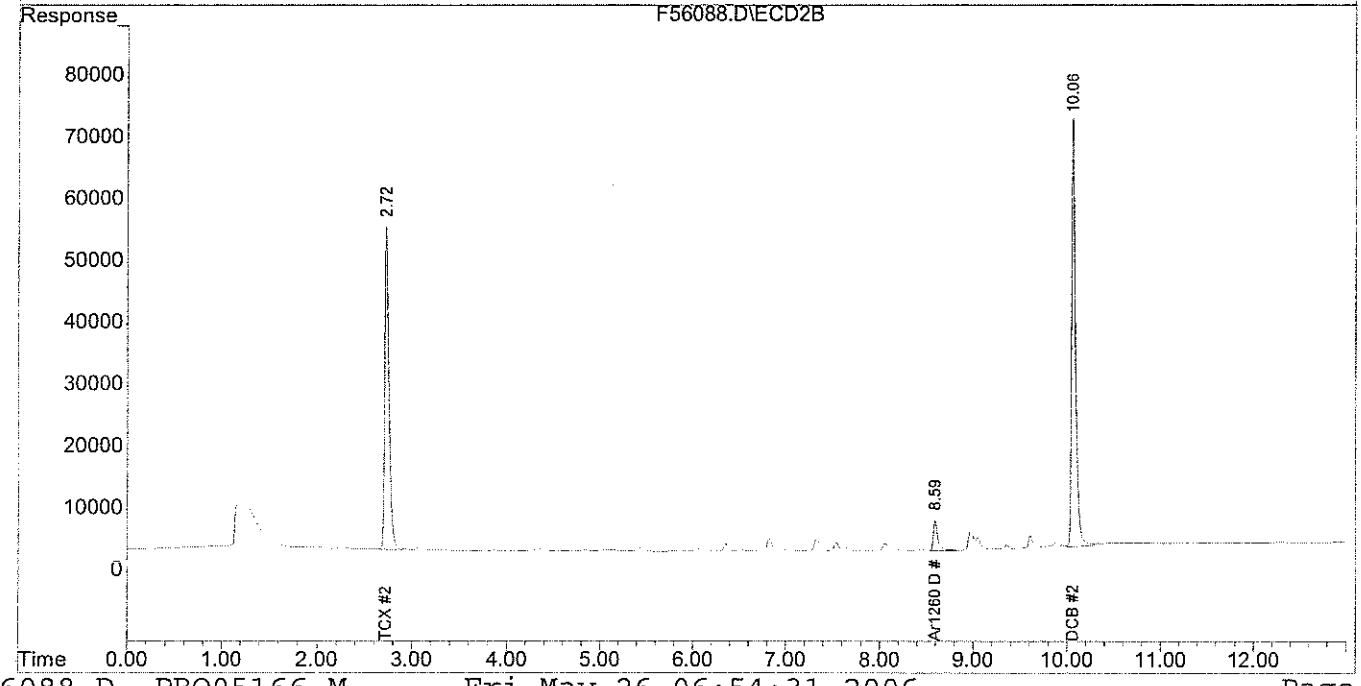
Quantitation Report (QT Reviewed)

Signal #1 : C:\HPCHEM\2\DATA\052506-F\F56088.D\ECD1A.CH Vial: 11
Signal #2 : C:\HPCHEM\2\DATA\052506-F\F56088.D\ECD2B.CH
Acq On : 25 May 2006 5:51 pm Operator:
Sample : 56617-30, A/C Inst : Instrumen
Misc : SOIL Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 26 6:52 2006 Quant Results File: PBQ05166.RES

Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 10:06:32 2006
Response via : Multiple Level Calibration
DataAcq Meth : PSTQUICK.M

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film





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May 26, 2006

CLIENT SAMPLE ID

Project Name: GP

Project Number: 218736

Field Sample ID: W-7

SAMPLE DATA

Lab Sample ID: 56617-31
Matrix: Wipe
Percent Solid: N/A
Dilution Factor: 1.0
Collection Date: 05/24/06
Lab Receipt Date: 05/25/06
Extraction Date: 05/25/06
Analysis Date: 05/25/06

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit ug/wipe	Results ug/wipe
PCB-1016	0.5	U
PCB-1221	0.5	U
PCB-1232	0.5	U
PCB-1242	0.5	U
PCB-1248	0.5	U
PCB-1254	0.5	U
PCB-1260	0.5	U

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	100	%
Decachlorobiphenyl	108	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8081/82.

COMMENTS:

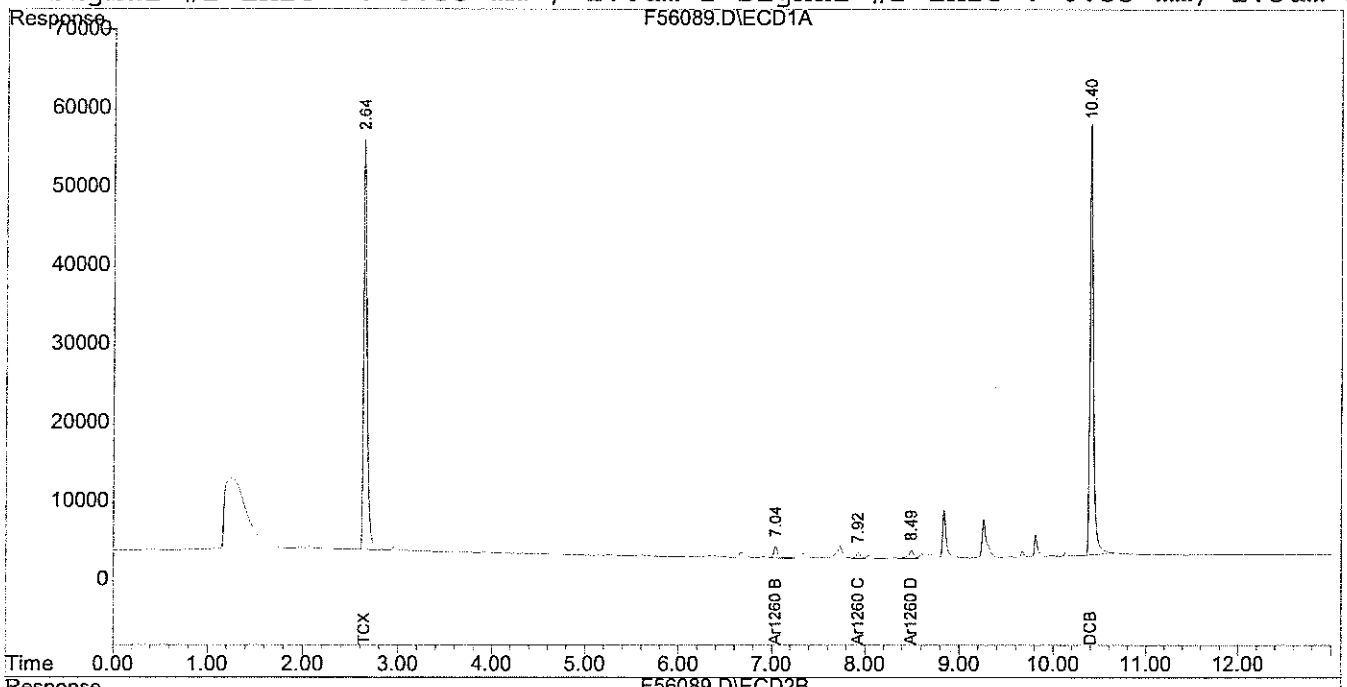
Authorized signature Milena Tull

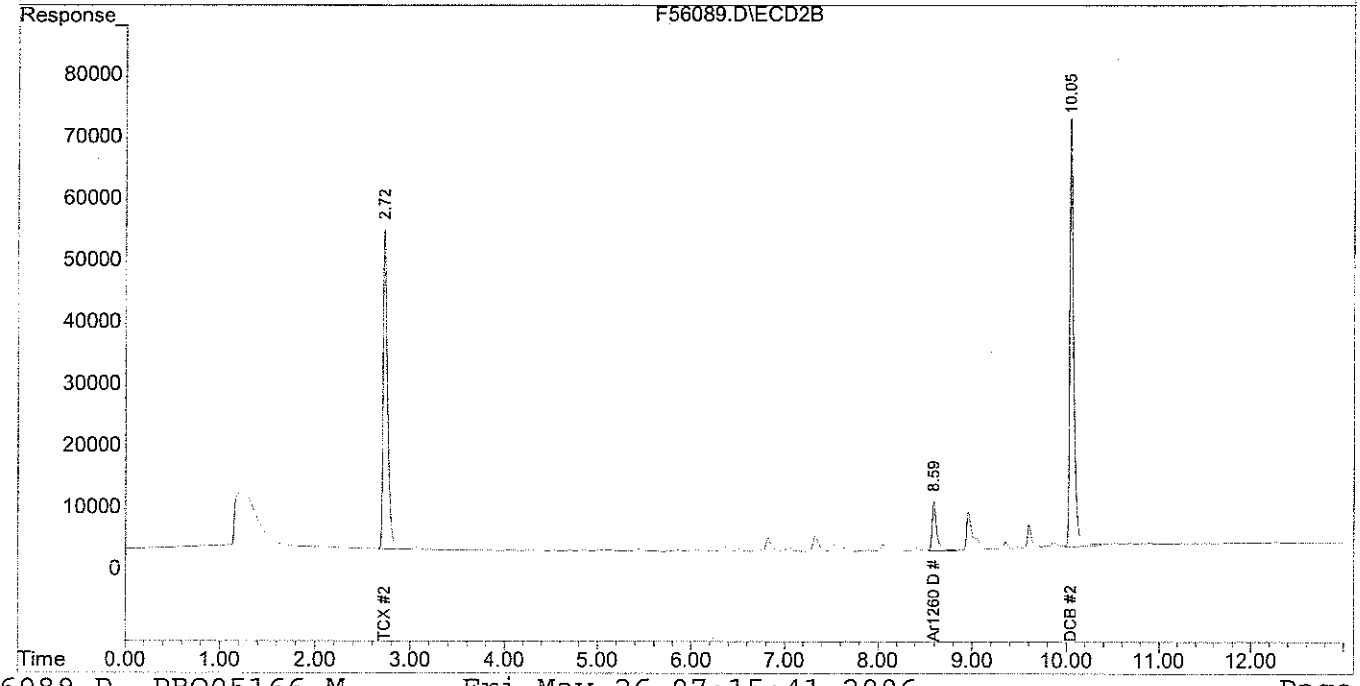
Quantitation Report (QT Reviewed)

Signal #1 : C:\HPCHEM\2\DATA\052506-F\F56089.D\ECD1A.CH Vial: 12
Signal #2 : C:\HPCHEM\2\DATA\052506-F\F56089.D\ECD2B.CH
Acq On : 25 May 2006 6:05 pm Operator:
Sample : 56617-31, A/C Inst : Instrumen
Misc : SOIL Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 26 7:15 2006 Quant Results File: PBQ05166.RES

Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 10:06:32 2006
Response via : Multiple Level Calibration
DataAcq Meth : PSTQUICK.M

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film





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May 26, 2006

CLIENT SAMPLE ID

Project Name: GP

Project Number: 218736

Field Sample ID: W-8

SAMPLE DATA

Lab Sample ID: 56617-32
Matrix: Wipe
Percent Solid: N/A
Dilution Factor: 1.0
Collection Date: 05/24/06
Lab Receipt Date: 05/25/06
Extraction Date: 05/25/06
Analysis Date: 05/25/06

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit µg/wipe	Results µg/wipe
PCB-1016	0.5	U
PCB-1221	0.5	U
PCB-1232	0.5	U
PCB-1242	0.5	U
PCB-1248	0.5	U
PCB-1254	0.5	U
PCB-1260	0.5	**5.0 D**

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	103	%
Decachlorobiphenyl	115	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8081/82.

Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3550B.

COMMENTS: D=Sample did not meet confirmation acceptance criteria for percent difference.

PCB Report

Authorized signature

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PCB CONFIRMATION RESULTS

SDG: 56617

Project No: Site ID: Location: ME

Instrument ID: F GC Columns: RTX-CLPesticidesI ID: 0.32(mm)
RTX-CLPesticidesII

Sample ID: 56617-32

COMPOUND	RESPONSE #1 ug/wipe	RESPONSE #2 ug/wipe	DIFFERENCE %	#	% DIFFERENCE Acceptance Limit
PCB 1260	3.0	5.0	50	*	40

Column to be used to flag % Difference with an asterisk

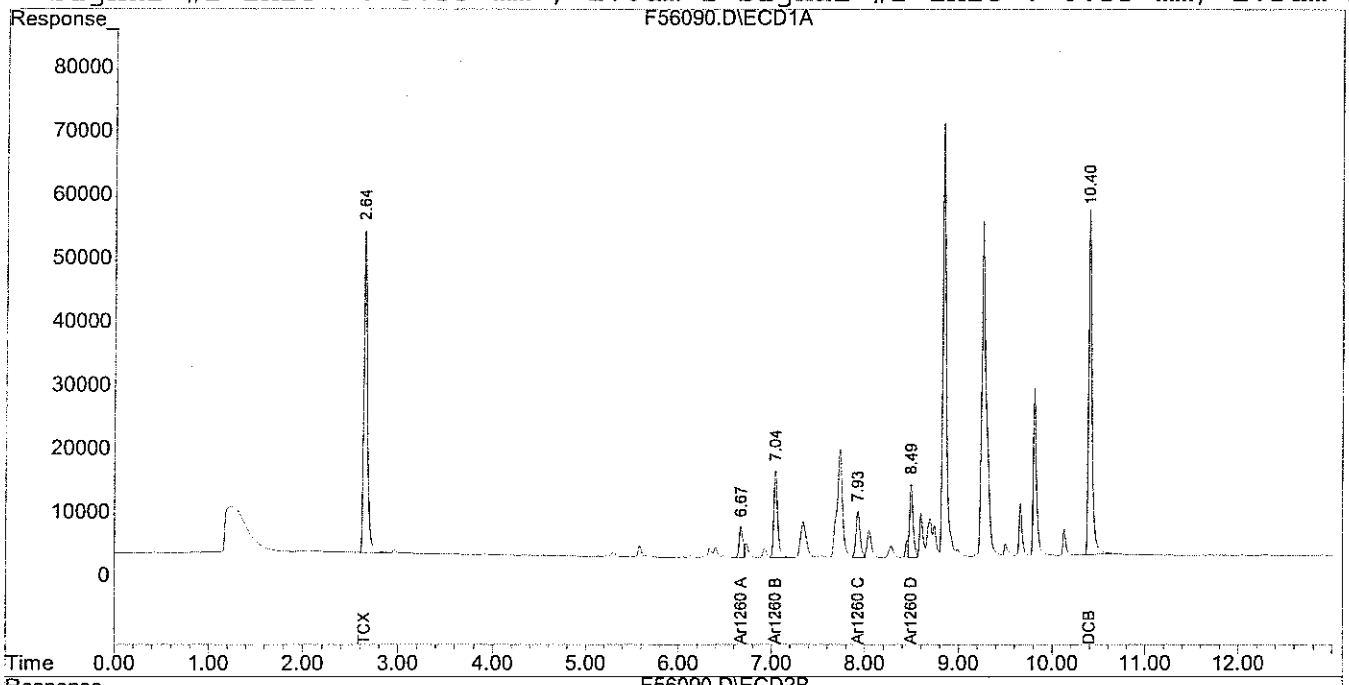
* Values outside of QC limits

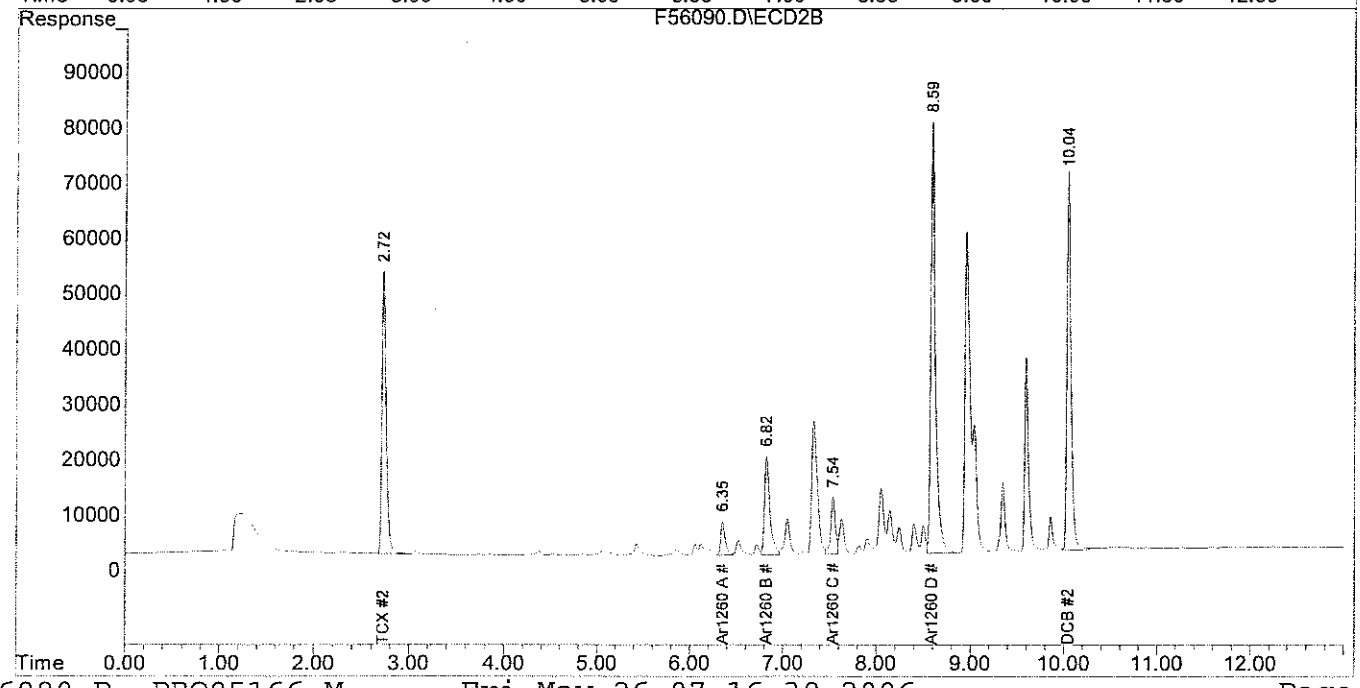
Quantitation Report (QT Reviewed)

Signal #1 : C:\HPCHEM\2\DATA\052506-F\F56090.D\ECD1A.CH Vial: 13
Signal #2 : C:\HPCHEM\2\DATA\052506-F\F56090.D\ECD2B.CH
Acq On : 25 May 2006 6:20 pm Operator:
Sample : 56617-32, A/C Inst : Instrumen
Misc : SOIL Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 26 7:16 2006 Quant Results File: PBQ05166.RES

Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 10:06:32 2006
Response via : Multiple Level Calibration
DataAcq Meth : PSTQUICK.M

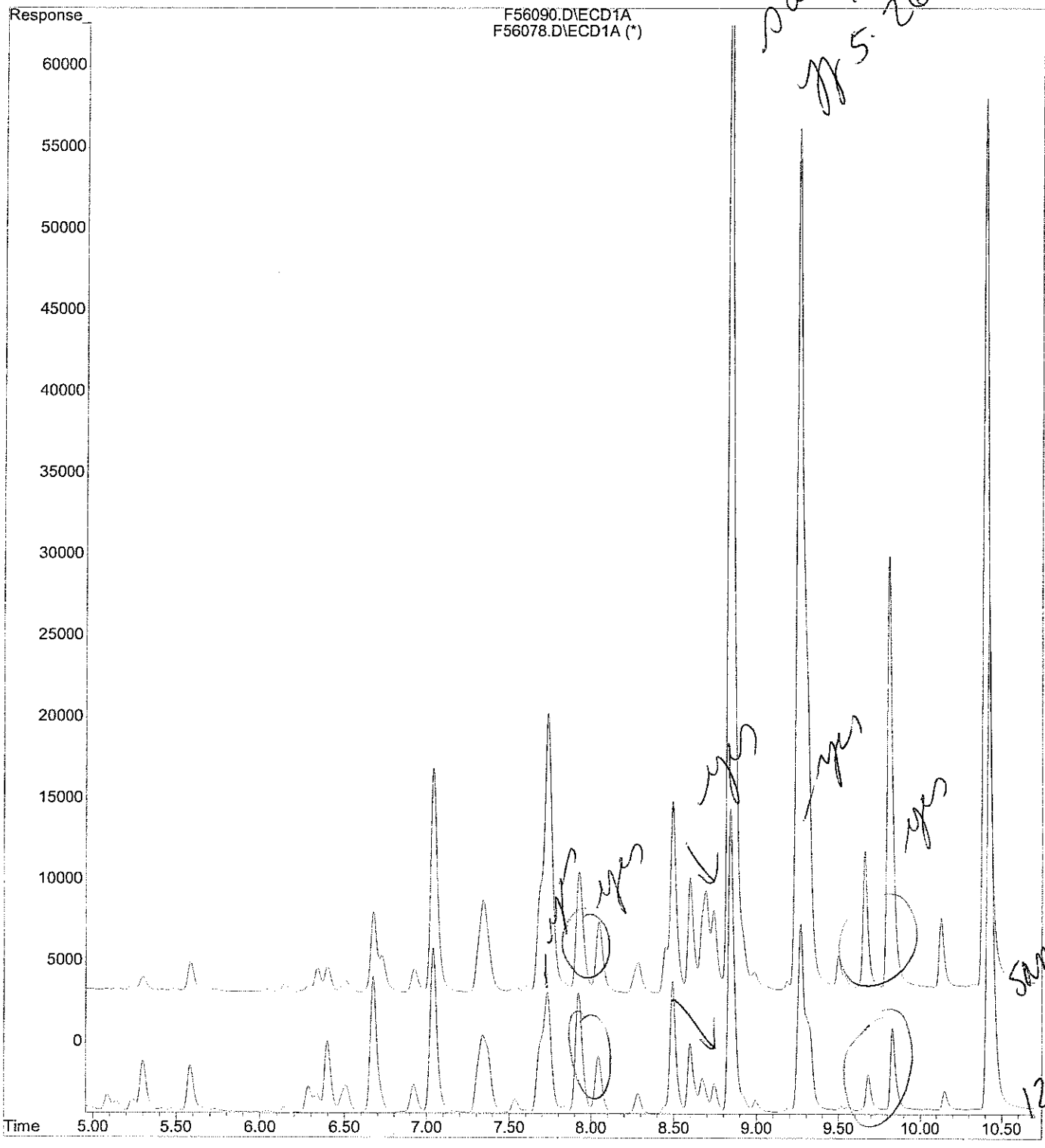
Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film





File : C:\HPCHEM\2\DATA\052506-F\F56090.D
Operator :
Acquired : 25 May 2006 6:20 pm using AcqMethod PSTQUICK.M
Instrument : Instrumen
Sample Name: 56617-32, A/C
Misc Info : SOIL
Vial Number: 13

Sample contains 1260 @ 5.26.06



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May 26, 2006

CLIENT SAMPLE ID

Project Name: GP

Project Number: 218736

Field Sample ID: W-9

SAMPLE DATA

Lab Sample ID: 56617-33
Matrix: Wipe
Percent Solid: N/A
Dilution Factor: 1.0
Collection Date: 05/24/06
Lab Receipt Date: 05/25/06
Extraction Date: 05/25/06
Analysis Date: 05/25/06

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit ug/wipe	Results ug/wipe
PCB-1016	0.5	U
PCB-1221	0.5	U
PCB-1232	0.5	U
PCB-1242	0.5	U
PCB-1248	0.5	U
PCB-1254	0.5	U
PCB-1260	0.5	**0.8**

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	95	%
Decachlorobiphenyl	108	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8081/82.

COMMENTS:

PCB Report

Authorized signature

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PCB CONFIRMATION RESULTS

SDG: 56617

Project No: Site ID: Location: ME

Instrument ID: F GC Columns: RTX-CLPesticidesI ID: 0.32(mm)
RTX-CLPesticidesII

Sample ID: 56617-33

COMPOUND	RESPONSE #1 ug/wipe	RESPONSE #2 ug/wipe	DIFFERENCE %	#	% DIFFERENCE Acceptance Limit
PCB 1260	0.8	0.7	18		40

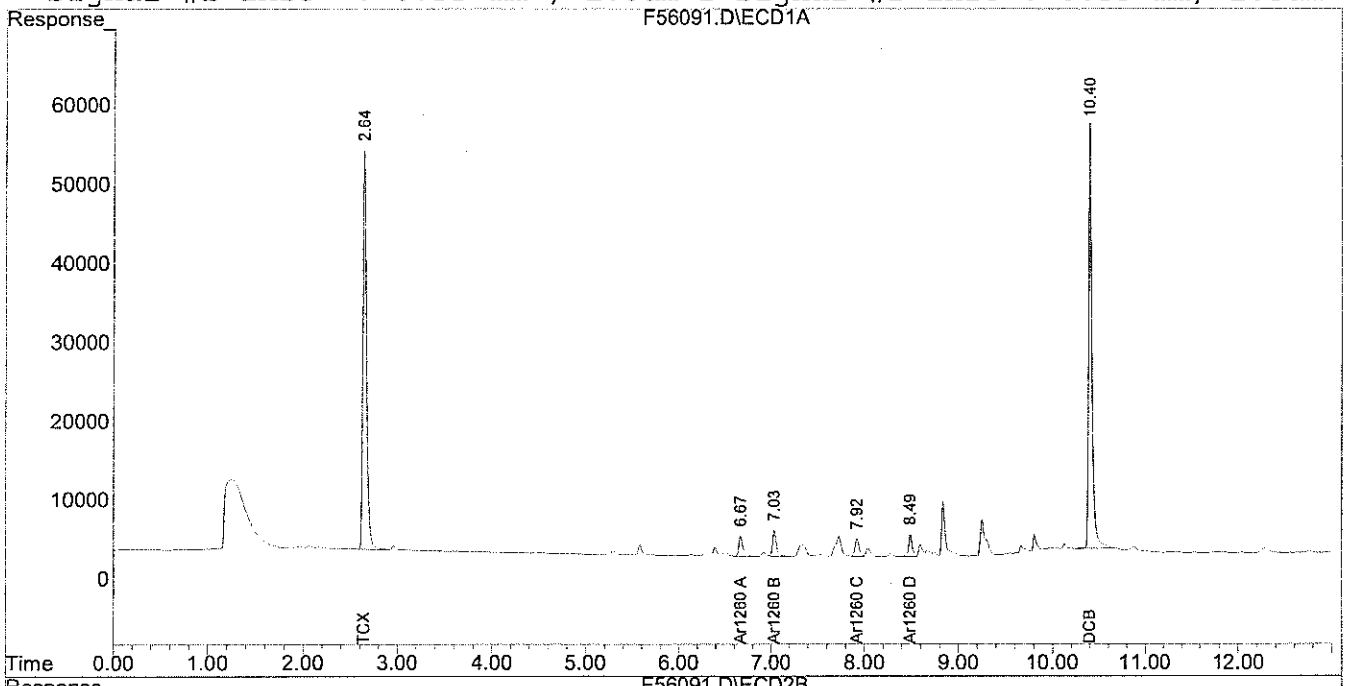
Column to be used to flag % Difference with an asterisk

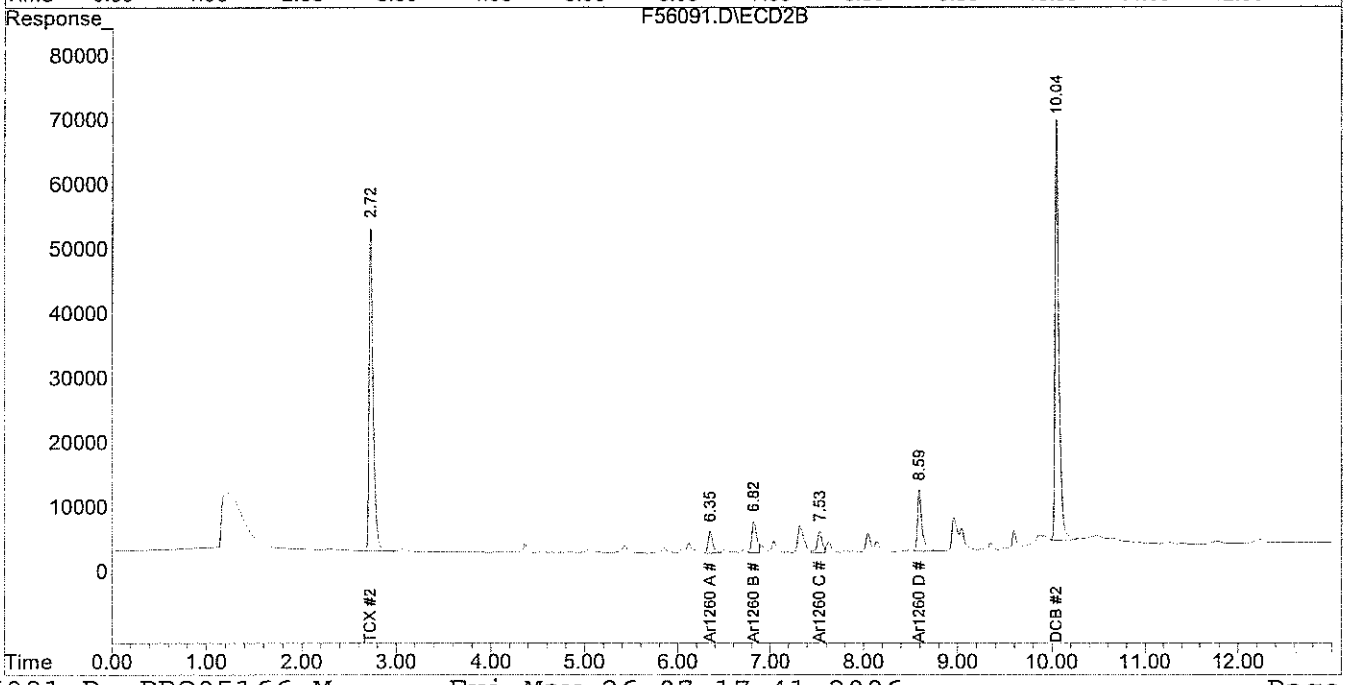
* Values outside of QC limits

Signal #1 : C:\HPCHEM\2\DATA\052506-F\F56091.D\ECD1A.CH Vial: 14
Signal #2 : C:\HPCHEM\2\DATA\052506-F\F56091.D\ECD2B.CH
Acq On : 25 May 2006 6:35 pm Operator:
Sample : 56617-33, A/C Inst : Instrumen
Misc : SOIL Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 26 7:17 2006 Quant Results File: PBQ05166.RES

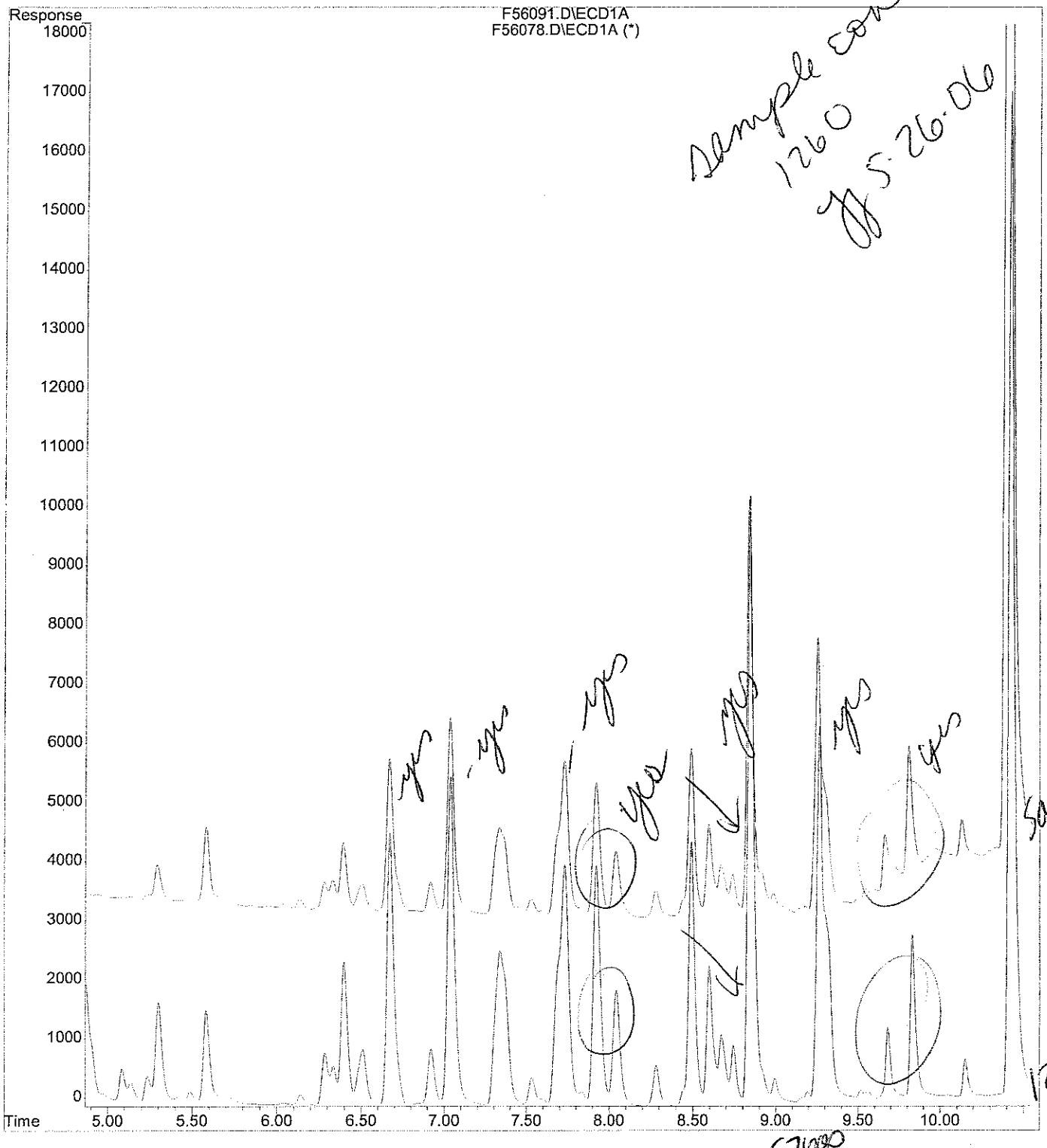
Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 10:06:32 2006
Response via : Multiple Level Calibration
DataAcq Meth : PSTQUICK.M

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film





File : C:\HPCHEM\2\DATA\052506-F\F56091.D
Operator :
Acquired : 25 May 2006 6:35 pm using AcqMethod PSTQUICK.M
Instrument : Instrumen
Sample Name: 56617-33, A/C
Misc Info : SOIL
Vial Number: 14



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Mr. Karl Kasper
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Portland ME 04102

May 26, 2006

SAMPLE DATA

Lab Sample ID:	56617-34
Matrix:	Wipe
Percent Solid:	N/A
Dilution Factor:	1.0
Collection Date:	05/24/06
Lab Receipt Date:	05/25/06
Extraction Date:	05/25/06
Analysis Date:	05/25/06

CLIENT SAMPLE ID

Project Name:	GP
Project Number:	218736
Field Sample ID:	W-10

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit ug/wipe	Results ug/wipe
PCB-1016	0.5	U
PCB-1221	0.5	U
PCB-1232	0.5	U
PCB-1242	0.5	U
PCB-1248	0.5	U
PCB-1254	0.5	U
PCB-1260	0.5	U

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	95	%
Decachlorobiphenyl	108	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8081/82.

COMMENTS:

PCB Report

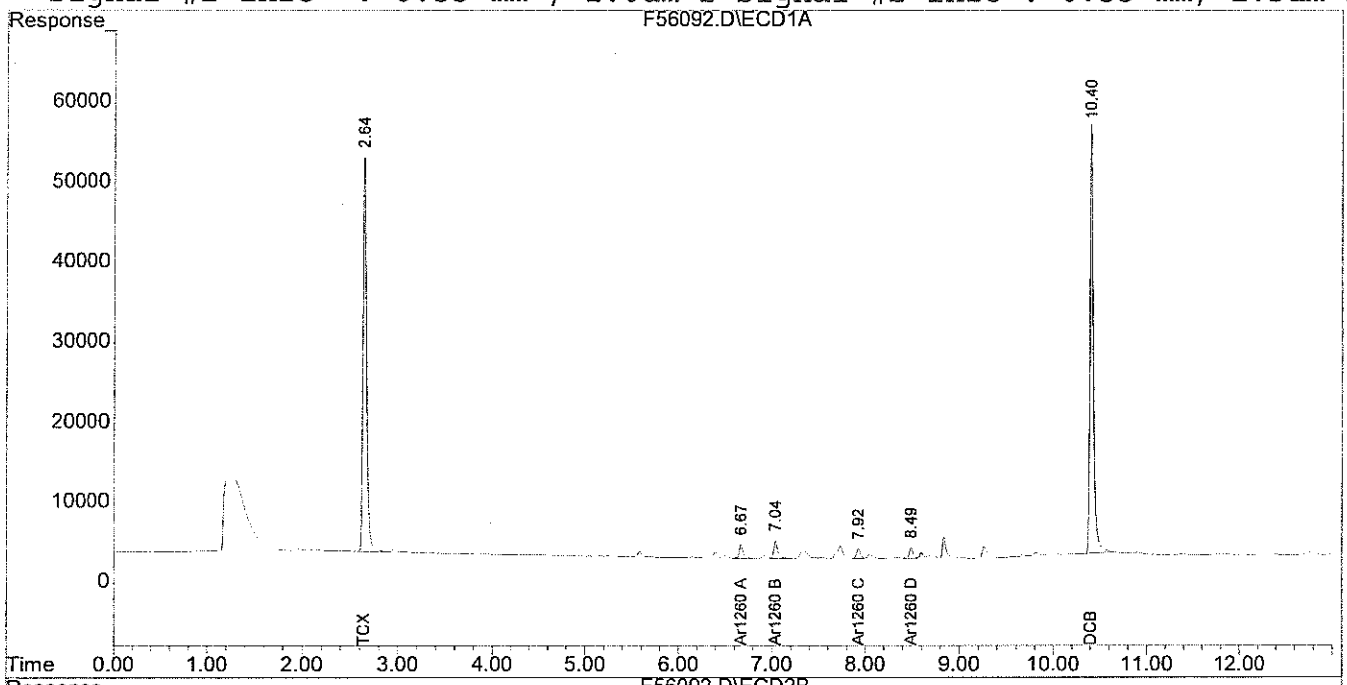
Authorized signature [signature]

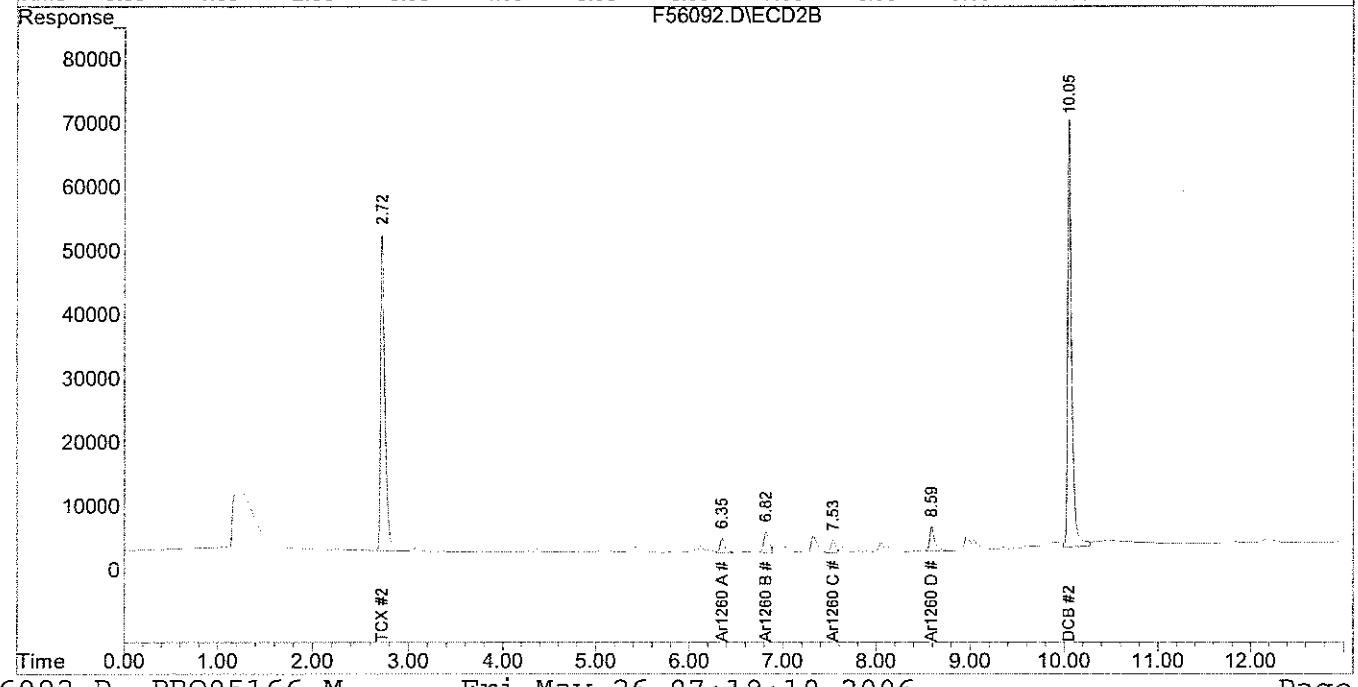
Quantitation Report (QT Reviewed)

Signal #1 : C:\HPCHEM\2\DATA\052506-F\F56092.D\ECD1A.CH Vial: 15
Signal #2 : C:\HPCHEM\2\DATA\052506-F\F56092.D\ECD2B.CH
Acq On : 25 May 2006 6:49 pm Operator:
Sample : 56617-34, A/C Inst : Instrumen
Misc : SOIL Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 26 7:19 2006 Quant Results File: PBQ05166.RES

Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 10:06:32 2006
Response via : Multiple Level Calibration
DataAcq Meth : PSTQUICK.M

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film





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Woodard & Curran
41 Hutchins Drive
Portland ME 04102

May 26, 2006

SAMPLE DATA

CLIENT SAMPLE ID

Project Name: GP

Project Number: 218736

Field Sample ID: W-11

Lab Sample ID: 56617-35
Matrix: Wipe
Percent Solid: N/A
Dilution Factor: 1.0
Collection Date: 05/24/06
Lab Receipt Date: 05/25/06
Extraction Date: 05/25/06
Analysis Date: 05/25/06

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit ug/wipe	Results ug/wipe
PCB-1016	0.5	U
PCB-1221	0.5	U
PCB-1232	0.5	U
PCB-1242	0.5	U
PCB-1248	0.5	U
PCB-1254	0.5	U
PCB-1260	0.5	U

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	98	%
Decachlorobiphenyl	108	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8081/82.

COMMENTS:

PCB Report

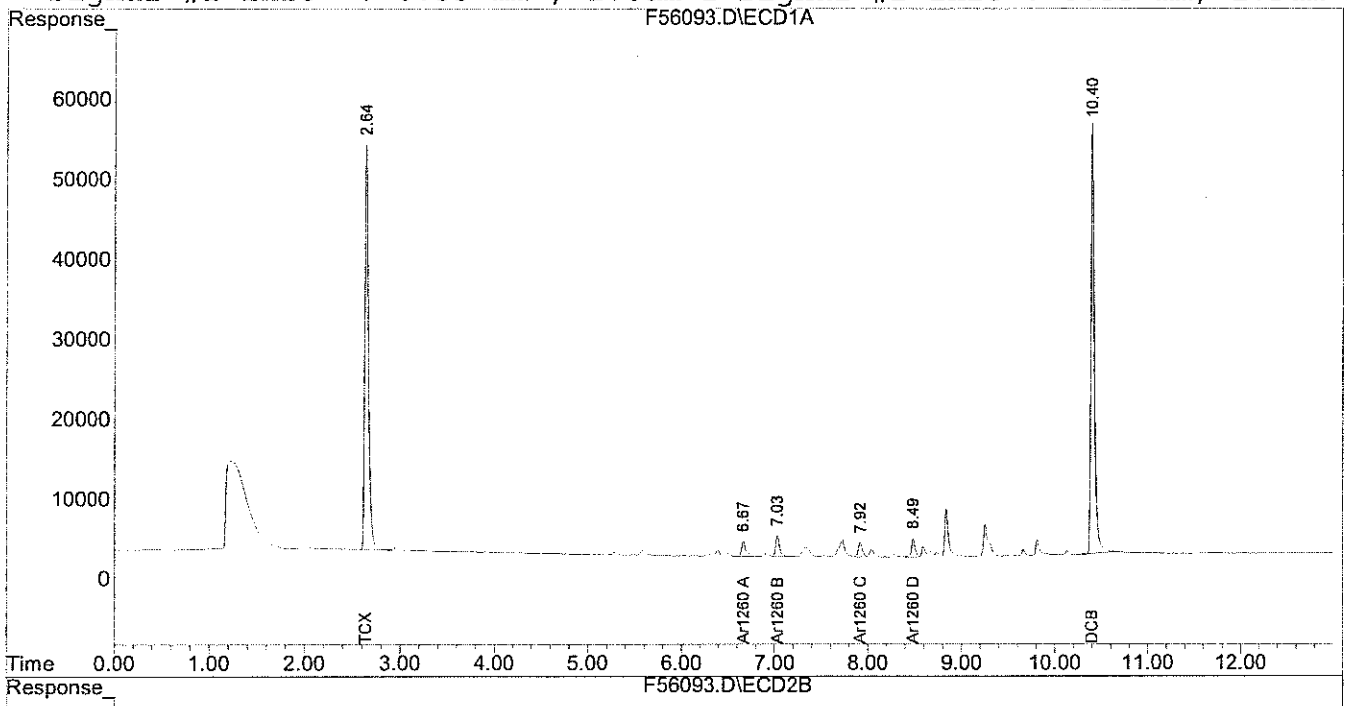
Authorized signature Melissa Hall

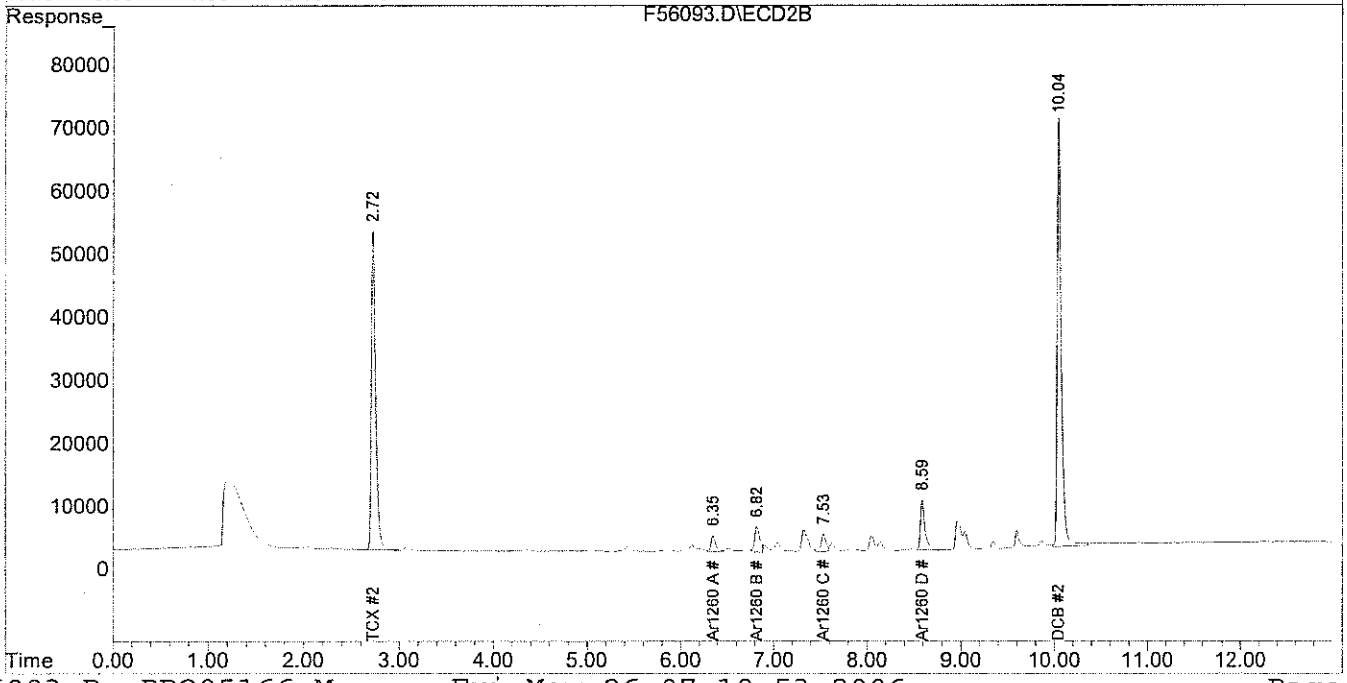
Quantitation Report (QT Reviewed)

Signal #1 : C:\HPCHEM\2\DATA\052506-F\F56093.D\ECD1A.CH Vial: 16
Signal #2 : C:\HPCHEM\2\DATA\052506-F\F56093.D\ECD2B.CH
Acq On : 25 May 2006 7:19 pm Operator:
Sample : 56617-35, A/C Inst : Instrumen
Misc : SOIL Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 26 7:19 2006 Quant Results File: PBQ05166.RES

Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 10:06:32 2006
Response via : Multiple Level Calibration
DataAcq Meth : PSTQUICK.M

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film





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41 Hutchins Drive
Portland ME 04102

May 26, 2006

SAMPLE DATA

CLIENT SAMPLE ID

Project Name: GP

Project Number: 218736

Field Sample ID: W-12

Lab Sample ID: 56617-36
Matrix: Wipe
Percent Solid: N/A
Dilution Factor: 1.0
Collection Date: 05/24/06
Lab Receipt Date: 05/25/06
Extraction Date: 05/25/06
Analysis Date: 05/25/06

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit ug/wipe	Results ug/wipe
PCB-1016	0.5	U
PCB-1221	0.5	U
PCB-1232	0.5	U
PCB-1242	0.5	U
PCB-1248	0.5	U
PCB-1254	0.5	U
PCB-1260	0.5	5.5

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	95	%
Decachlorobiphenyl	108	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8081/82.

COMMENTS:

PCB Report

Authorized signature Melena Hall

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PCB CONFIRMATION RESULTS

SDG: 56617

Project No: Site ID: Location: ME

Instrument ID: F GC Columns: RTX-CLPesticidesI ID: 0.32(mm)
RTX-CLPesticidesII

Sample ID: 56617-36

COMPOUND	RESPONSE #1 ug/wipe	RESPONSE #2 ug/wipe	DIFFERENCE %	#	% DIFFERENCE Acceptance Limit
PCB 1260	5.5	6.9	23		40

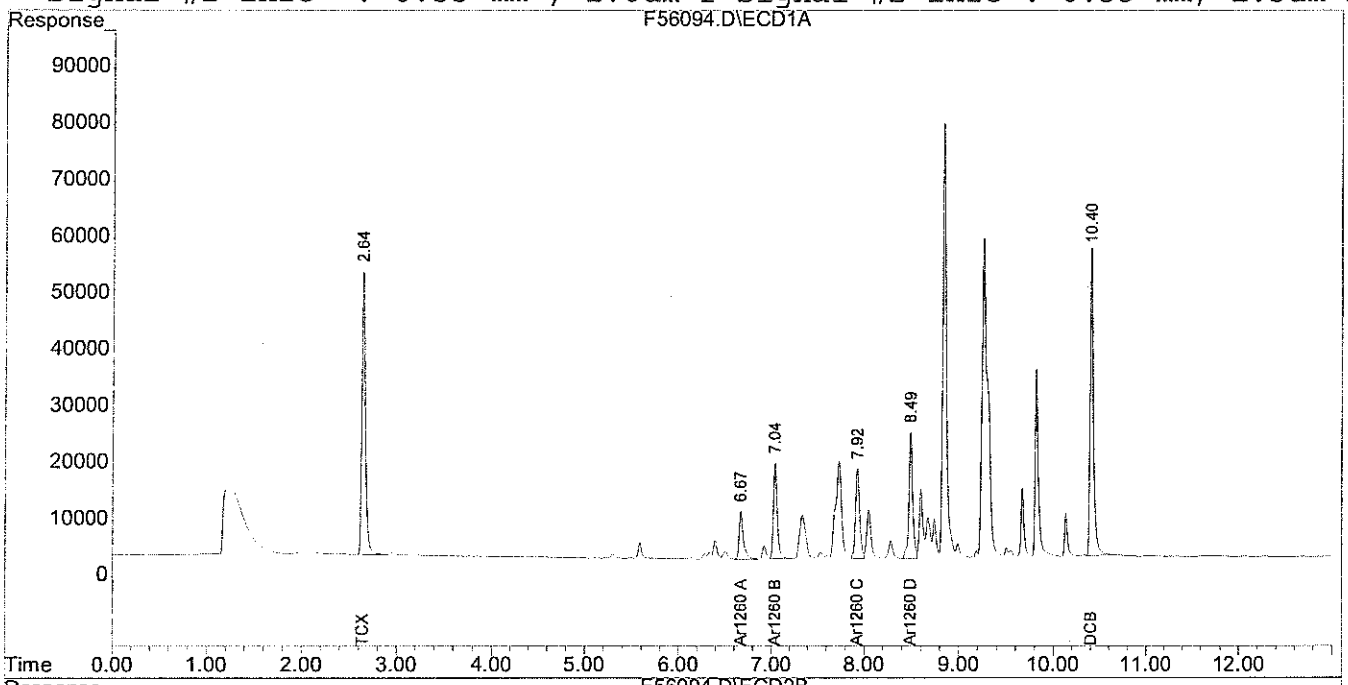
Column to be used to flag % Difference with an asterisk

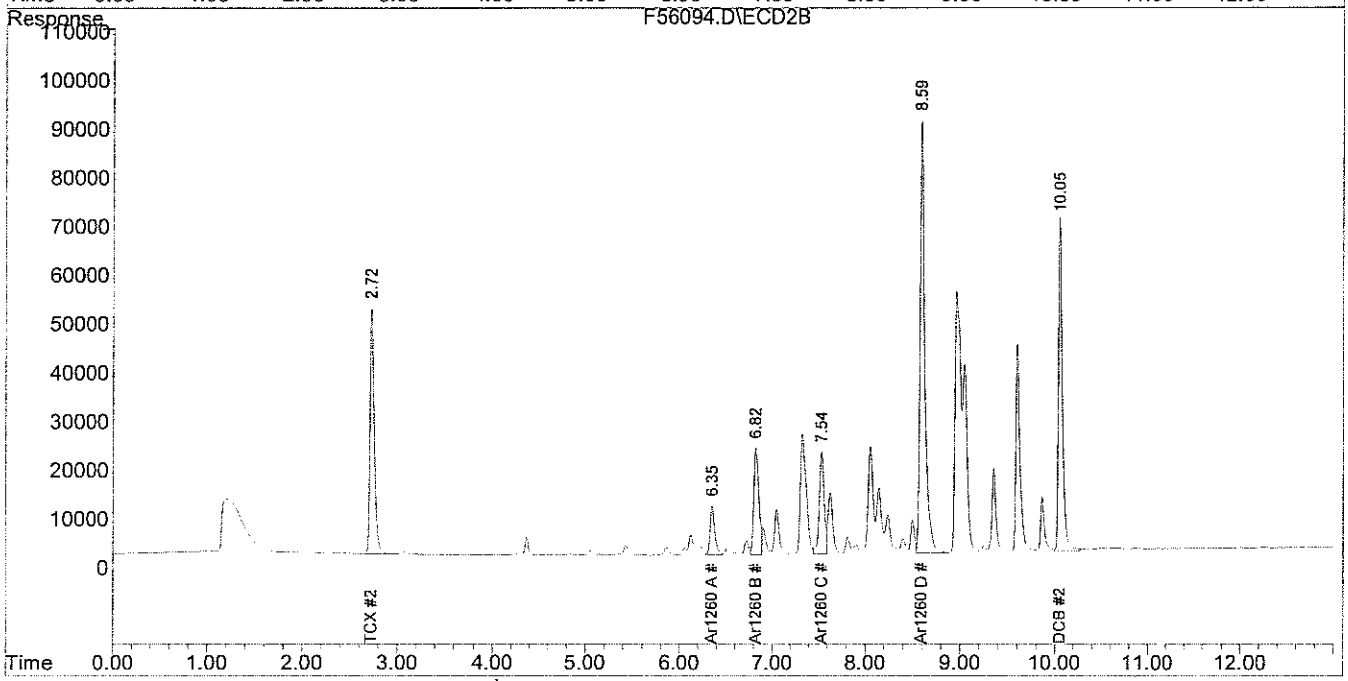
* Values outside of QC limits

Signal #1 : C:\HPCHEM\2\DATA\052506-F\F56094.D\ECD1A.CH Vial: 17
Signal #2 : C:\HPCHEM\2\DATA\052506-F\F56094.D\ECD2B.CH
Acq On : 25 May 2006 7:33 pm Operator:
Sample : 56617-36, A/C Inst : Instrumen
Misc : SOIL Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 26 7:20 2006 Quant Results File: PBQ05166.RES

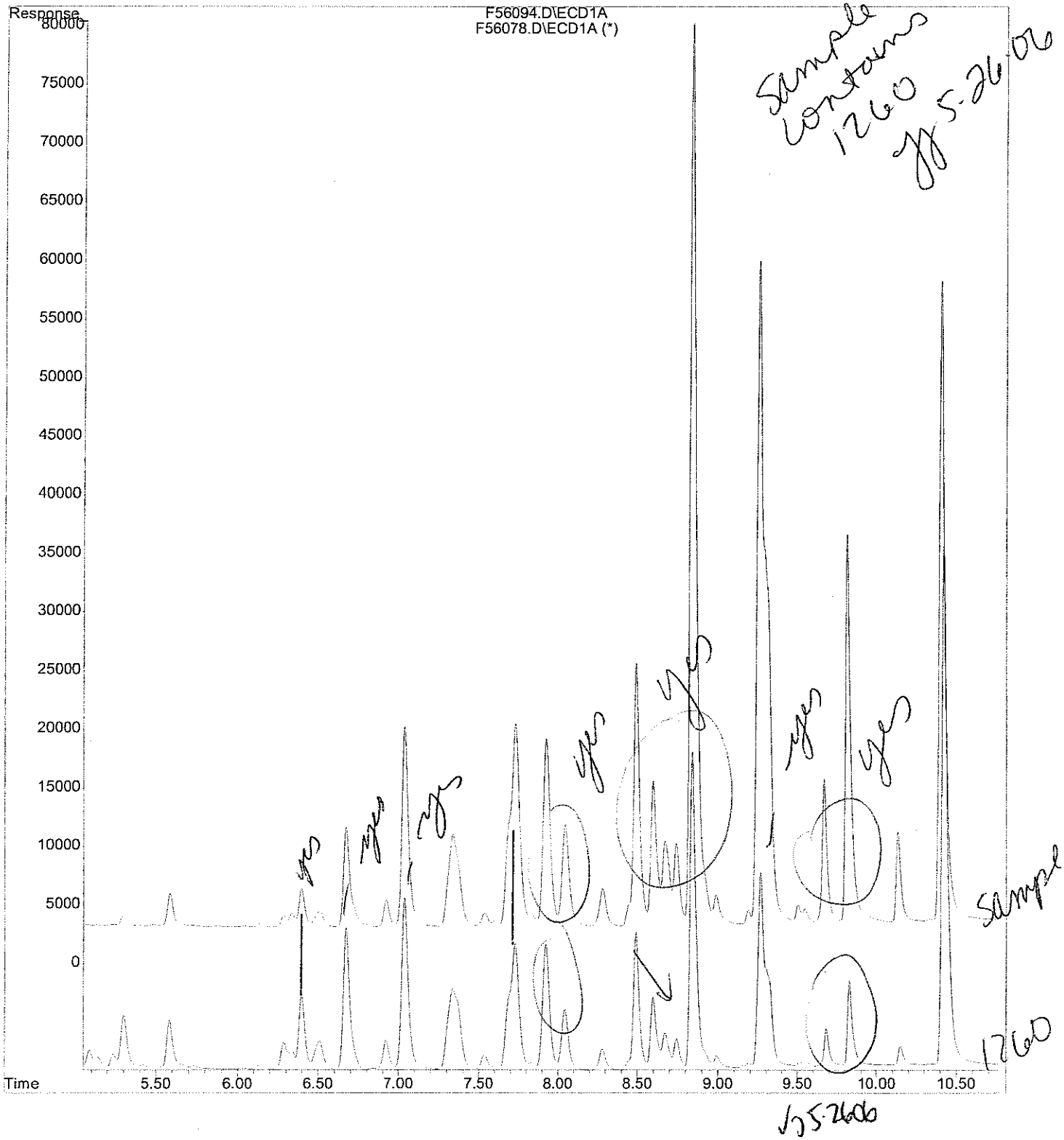
Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 10:06:32 2006
Response via : Multiple Level Calibration
DataAcq Meth : PSTQUICK.M

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film





File : C:\HPCHEM\2\DATA\052506-F\F56094.D
Operator :
Acquired : 25 May 2006 7:33 pm using AcqMethod PSTQUICK.M
Instrument : Instrumen
Sample Name: 56617-36, A/C
Misc Info : SOIL
Vial Number: 17



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May 26, 2006

CLIENT SAMPLE ID

Project Name: GP

Project Number: 218736

Field Sample ID: W-13

SAMPLE DATA

Lab Sample ID: 56617-37
Matrix: Wipe
Percent Solid: N/A
Dilution Factor: 1.0
Collection Date: 05/24/06
Lab Receipt Date: 05/25/06
Extraction Date: 05/25/06
Analysis Date: 05/25/06

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit ug/wipe	Results ug/wipe
PCB-1016	0.5	U
PCB-1221	0.5	U
PCB-1232	0.5	U
PCB-1242	0.5	U
PCB-1248	0.5	U
PCB-1254	0.5	U
PCB-1260	0.5	**0.9**

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	98	%
Decachlorobiphenyl	110	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8081/82.

COMMENTS:

PCB Report

Authorized signature

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PCB CONFIRMATION RESULTS

SDG: 56617

Project No: Site ID: Location: ME

Instrument ID: F GC Columns: RTX-CLPesticidesI ID: 0.32(mm)
RTX-CLPesticidesII

Sample ID: 56617-37

COMPOUND	RESPONSE #1 ug/wipe	RESPONSE #2 ug/wipe	DIFFERENCE %	#	% DIFFERENCE Acceptance Limit
PCB 1260	0.9	0.8	15		40

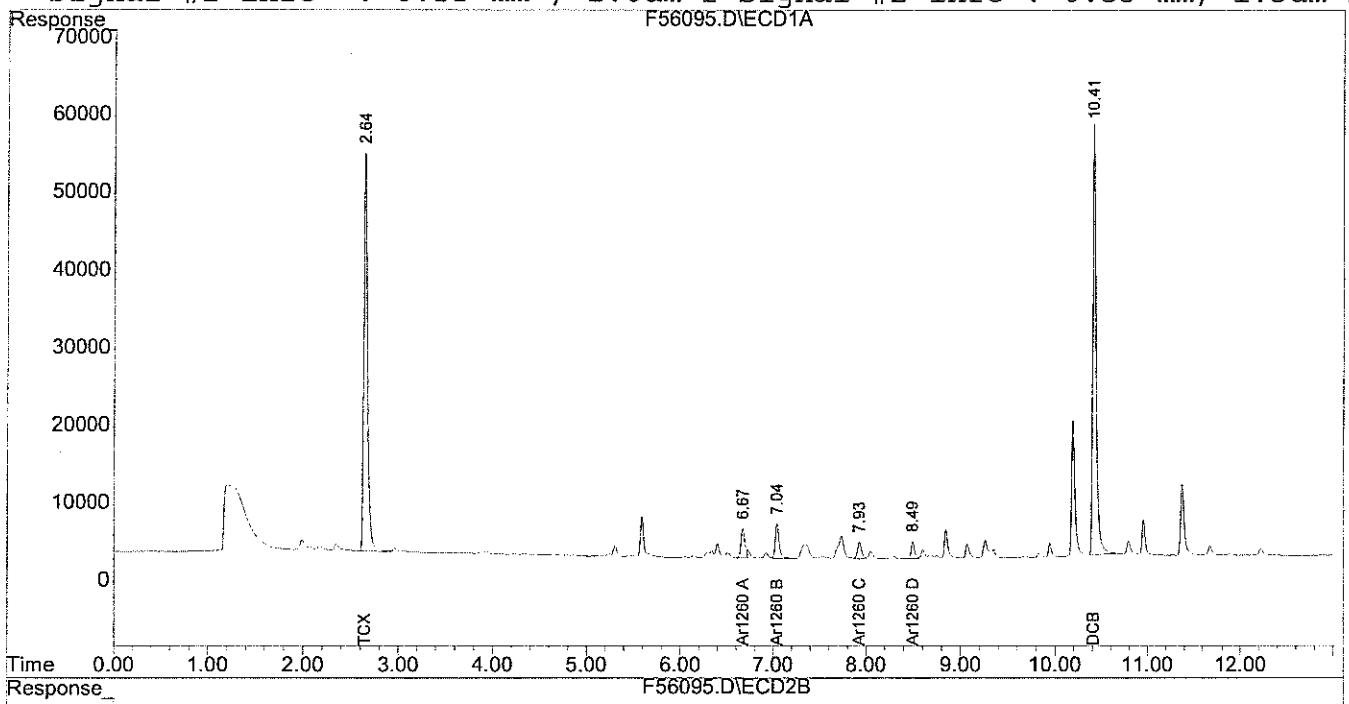
Column to be used to flag % Difference with an asterisk

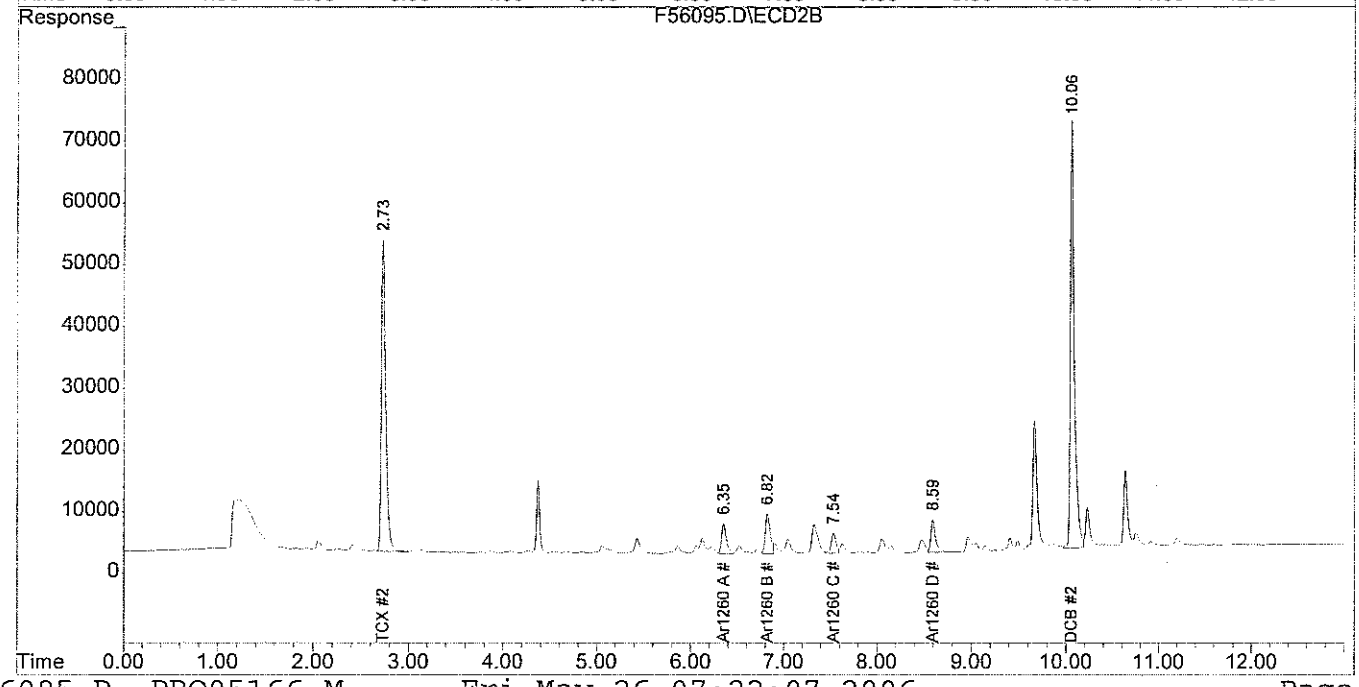
* Values outside of QC limits

Signal #1 : C:\HPCHEM\2\DATA\052506-F\F56095.D\ECD1A.CH Vial: 18
Signal #2 : C:\HPCHEM\2\DATA\052506-F\F56095.D\ECD2B.CH
Acq On : 25 May 2006 7:48 pm Operator:
Sample : 56617-37, A/C Inst : Instrumen
Misc : SOIL Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 26 7:21 2006 Quant Results File: PBQ05166.RES

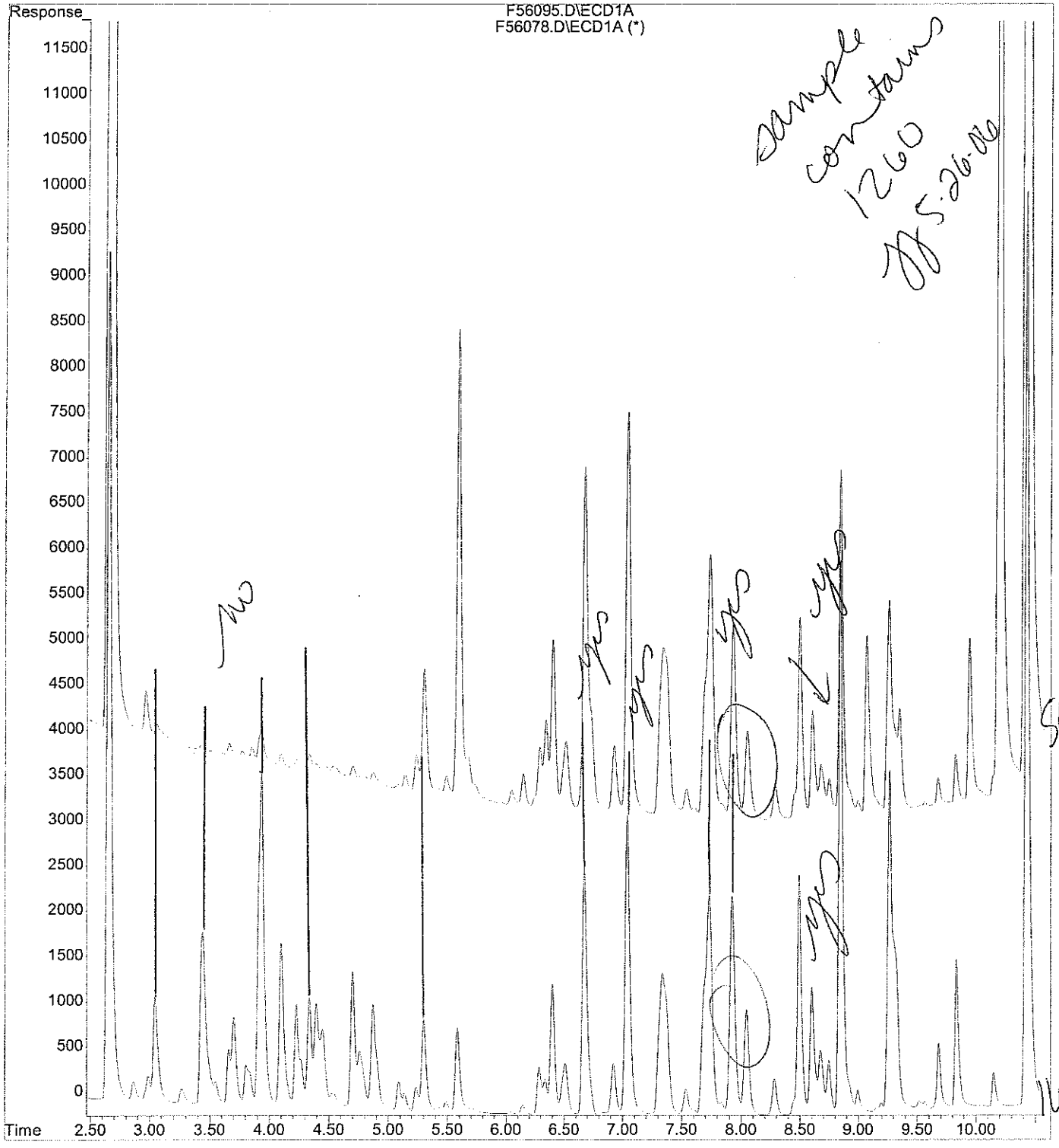
Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 10:06:32 2006
Response via : Multiple Level Calibration
DataAcq Meth : PSTQUICK.M

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film





File : C:\HPCHEM\2\DATA\052506-F\F56095.D
Operator :
Acquired : 25 May 2006 7:48 pm using AcqMethod PSTQUICK.M
Instrument : Instrumen
Sample Name: 56617-37, A/C
Misc Info : SOIL
Vial Number: 18



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PCB
QC FORMS

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PCB SOLID
SYSTEM MONITORING COMPOUND RECOVERY

Project #: 218736 Quote #: SDG#: 56617

Site ID: GP Location: ME

ıstrument ID: F
GC Column: RTX-CLPesticides I ID: 0.32(mm)
RTX-CLPesticides II

		COLUMN #1			COLUMN #2			TOTAL	
	SAMPLE ID	SMC1	#	SMC2	#	SMC1	#	SMC2	#	OUT	#
01	P4016 16/60 0.01	NA		NA		NA		NA		0	
02	P4017 16/60 0.02	NA		NA		NA		NA		0	
03	P4018 16/60 0.05	NA		NA		NA		NA		0	
04	P4019 16/60 0.1	NA		NA		NA		NA		0	
05	P4020 16/60 0.25	116		109		116		121		0	
06	P4021 16/60 0.5	112		104		110		108		0	
07	P4022 16/60 1.0	NA		NA		NA		NA		0	
08	P4024 16/60 0.5 C	104		104		105		107		0	
09	P3883 AR 1221	NA		NA		NA		NA		0	
10	P3884 AR 1232	NA		NA		NA		NA		0	
11	P3885 AR 1242	NA		NA		NA		NA		0	
12	P3886 AR 1248	NA		NA		NA		NA		0	
13	P3994 AR 1254	NA		NA		NA		NA		0	
14	P3887 AR 1262	NA		NA		NA		NA		0	
15	P3888 AR 1268	NA		NA		NA		NA		0	
16	P4057 AR 1254 0.	NA		NA		NA		NA		0	
17	P4056 AR 1254 0.	NA		NA		NA		NA		0	
18	P4055 AR 1254 0.	NA		NA		NA		NA		0	
19	P4054 AR 1254 0.	NA		NA		NA		NA		0	
20	P4053 AR 1254 0.	NA		NA		NA		NA		0	
21	P4052 AR 1254 1.	NA		NA		NA		NA		0	
22	P4010 AR 1254 0.	102		103		105		108		0	
23	ZZZZ	81		17	*	87		14	*	2	
24	ZZZZ	85		21	*	95		11	*	2	
25	ZZZZ	87		15	*	95		16	*	2	
26	ZZZZ	86		14	*	94		15	*	2	
27	ZZZZ	77		81		80		81		0	
28	ZZZZ	62		33	*	87		33	*	2	
29	ZZZZ	83		10	*	96		9	*	2	
30	P4021 16/60 0.5	99		103		102		109		0	
31	P4053 AR 1254 0.	NA		NA		NA		NA		0	
32											
33											
34											
35											
36											
37											
38											
39											
40											
41											
42											

QC LIMITS
SMC1 = TCX 46-123
SMC2 = DCB 38-123
*=Values outside of protocol required QC limits
D=System Monitoring Compound diluted out

analytics environmental laboratory LLC

PCB SOLID
SYSTEM MONITORING COMPOUND RECOVERY

Project #: 218736 Quote #: SDG#: 56617

Site ID: GP Location: ME

ıstrument ID: F

GC Column: RTX-CLPesticides I ID: 0.32(mm)

RTX-CLPesticides II

		COLUMN #1				COLUMN #2				TOTAL	
	SAMPLE ID	SMC1	#	SMC2	#	SMC1	#	SMC2	#	OUT	#
01	P4020 16/60 0.25	98		100		104		108		0	
02	B05246PAS, A/C	84		68		99		103		0	
03	L05246PASB, A/C	98		86		107		103		0	
04	LD05246PASB, A	100		90		108		103		0	
05	ZZZZZ	95		87		101		99		0	
06	ZZZZZ	91		78		97		88		0	
07	ZZZZZ	94		79		98		94		0	
08	ZZZZZ	94		80		99		86		0	
09	ZZZZZ	81		77		85		72		0	
10	ZZZZZ	86		240	*	90		204	*	2	
11	56606-6, MS, A/C	86		126	*	91		128	*	2	
12	56606-6, MSD, A/	87		139	*	93		144	*	2	
13	P4021 16/60 0.5	98		97		104		97		0	
14											
15											
16											
17											
18											
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39											
40											
41											
42											

QC LIMITS

SMC1 = TCX 46-123

SMC2 = DCB 38-123

*=Values outside of protocol required QC limits

D=System Monitoring Compound diluted out

analytics environmental laboratory LLC

PCB SOLID
SYSTEM MONITORING COMPOUND RECOVERY

Project #: 218736 Quote #: SDG#: 56617

Site ID: GP Location: ME

ıstrument ID: F

GC Column: RTX-CLPesticides I ID: 0.32(mm)
RTX-CLPesticides II

		COLUMN #1			COLUMN #2			TOTAL	
	SAMPLE ID	SMC1	#	SMC2	#	SMC1	#	SMC2	#	OUT	#
01	P4020 16/60 0.25	96		106		102		108		0	
02	B05256PWIPE, A	92		101		97		110		0	
03	L05256PWIPE, A	102		110		107		116		0	
04	LD05256PWIPE,	99		108		105		115		0	
05	56617-24, A/C	92		103		97		112		0	
06	56617-25, A/C	97		107		103		112		0	
07	56617-26, A/C	97		107		103		113		0	
08	56617-27, A/C	96		105		103		115		0	
09	ZZZZZ	97		109		103		122		0	
10	56617-29, A/C	98		110		105		118		0	
11	56617-30, A/C	99		109		105		117		0	
12	56617-31, A/C	99		109		104		117		0	
13	56617-32, A/C	96		108		102		114		0	
14	56617-33, A/C	96		109		101		108		0	
15	56617-34, A/C	94		107		100		115		0	
16	56617-35, A/C	96		106		102		115		0	
17	56617-36, A/C	96		108		101		116		0	
18	56617-37, A/C	98		111		103		120		0	
19	P4021 16/60 0.5	100		107		102		109		0	
20											
21											
22											
23											
24											
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38											
39											
40											
41											
42											

QC LIMITS
SMC1 = TCX 46-123
SMC2 = DCB 38-123
*=Values outside of protocol required QC limits
D=System Monitoring Compound diluted out

analytics environmental laboratory LLC

PCB SOLID
SYSTEM MONITORING COMPOUND RECOVERY

Project #: 218736 Quote #: SDG#: 56617

Site ID: GP Location: ME

ɔstrument ID: F

GC Column: RTX-CLPesticides I ID: 0.32(mm)
RTX-CLPesticides II

		COLUMN #1		COLUMN #2		TOTAL	
	SAMPLE ID	SMC1 #	SMC2 #	SMC1 #	SMC2 #	OUT	#
01	P4021 16/60 0.5	100	107	102	109	0	
02	B05256PAS, A/C	99	109	103	114	0	
03	L05256PAS, A/C	101	108	104	110	0	
04	56617-1, A/C	91	96	95	104	0	
05	56617-2, A/C	91	86	94	92	0	
06	56617-6, A/C	78	83	83	73	0	
07	56617-7, A/C	66	85	71	89	0	
08	56617-8, A/C	69	79	73	73	0	
09	56617-9, A/C	84	86	87	115	0	
10	56617-10, A/C	75	86	80	77	0	
11	56617-11, A/C	79	75	83	84	0	
12	56617-12, A/C	92	101	96	104	0	
13	56617-13, A/C	92	100	96	104	0	
14	P4021 16/60 0.5	100	108	102	109	0	
15							
16							
17							
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19							
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35							
36							
37							
38							
39							
40							
41							
42							

QC LIMITS
SMC1 = TCX 46-123
SMC2 = DCB 38-123
*=Values outside of protocol required QC limits
D=System Monitoring Compound diluted out

analytics environmental laboratory LLC

PCB SOLID
SYSTEM MONITORING COMPOUND RECOVERY

Project #: 218736 Quote #: SDG#: 56617

Site ID: GP Location: ME

ıstrument ID: F
GC Column: RTX-CLPesticides I ID: 0.32(mm)
RTX-CLPesticides II

		COLUMN #1			COLUMN #2			TOTAL	
	SAMPLE ID	SMC1	#	SMC2	#	SMC1	#	SMC2	#	OUT	#
01	P4021 16/60 0.5	100		108		102		109		0	
02	ZZZZZ	92		147	*	96		189	*	0	
03	ZZZZZ	D		D		D		D		0	
04	56617-4, 1:10, A/	90		90		80		80		0	
05	56617-5, A/C	99		78		110		85		0	
06	P4021 16/60 0.5	98		93		101		92		0	
07											
08											
09											
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QC LIMITS
SMC1 = TCX 46-123
SMC2 = DCB 38-123
*=Values outside of protocol required QC limits
D=System Monitoring Compound diluted out

analytics environmental laboratory LLC

PCB SOLID
SYSTEM MONITORING COMPOUND RECOVERY

Project #: 218736 Quote #: SDG#: 56617

Site ID: GP Location: ME

ıstrument ID: F

GC Column: RTX-CLPesticides I ID: 0.32(mm)
RTX-CLPesticides II

		COLUMN #1		COLUMN #2		TOTAL	
	SAMPLE ID	SMC1 #	SMC2 #	SMC1 #	SMC2 #	OUT	#
01	P4021 16/60 0.5	98	93	101	92	0	
02	ZZZZZ	NA	NA	NA	NA	0	
03	B05266PAS, A/C	102	99	108	114	0	
04	L05266PAS, A/C	97	97	103	105	0	
05	LD05266PAS, A/C	105	103	109	108	0	
06	56617-15, A/C	85	75	90	80	0	
07	ZZZZZ	86	115	92	81	0	
08	ZZZZZ	97	93	102	96	0	
09	56617-18, A/C	102	97	109	100	0	
10	56617-19, A/C	96	95	104	101	0	
11	56617-20, A/C	100	95	103	97	0	
12	56617-21, A/C	85	82	91	88	0	
13	56617-22, A/C	92	89	96	95	0	
14	56617-23, A/C	90	84	96	89	0	
15	56617-23, MS, A/C	88	77	93	80	0	
16	56617-23, MSD, A	91	78	99	81	0	
17	P4021 16/60 0.5	105	88	105	85	0	
18							
19							
20							
21							
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42							

QC LIMITS
SMC1 = TCX 46-123
SMC2 = DCB 38-123
*=Values outside of protocol required QC limits
D=System Monitoring Compound diluted out

analytics environmental laboratory LLC

PCB SOLID
SYSTEM MONITORING COMPOUND RECOVERY

Project #: 218736 Quote #: SDG#: 56617

Site ID: GP Location: ME

ıstrument ID: F
GC Column: RTX-CLPesticides I ID: 0.32(mm)
RTX-CLPesticides II

		COLUMN #1			COLUMN #2			TOTAL	
	SAMPLE ID	SMC1	#	SMC2	#	SMC1	#	SMC2	#	OUT	#
01	ZZZZZ	108		114		108		116		0	
02	P4020 16/60 0.25	106		114		108		114		0	
03	B05266PAS, RR A	109		112		109		120		0	
04	56617-28, 1:10, A/	D		D		D		D		0	
05	56617-14, 1:2000,	D		D		D		D		0	
06	56617-3, 1:5, A/C	90		90		95		95		0	
07	56617-16, 1:10, A/	90		120		130	*	70		0	
08	56617-17, 1:50, A/	D		D		D		D		0	
09	P4021 16/60 0.5	103		106		102		103		0	
10											
11											
12											
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42											

QC LIMITS
SMC1 = TCX 46-123
SMC2 = DCB 38-123
*=Values outside of protocol required QC limits
D=System Monitoring Compound diluted out

analytics environmental laboratory LLC

PCB SOLID
LABORATORY CONTROL SPIKE/ DUPLICATE
PERCENT RECOVERY

Project #: 218736 Quote #: SDG#: 56617

Site ID: Location: ME

Instrument ID: F

GC Column: RTX-CLPesticides I ID: 0.32(mm)

RTX-CLPesticides II

Laboratory Control Spike/Duplicate Sample ID: L05246PASB A/C/ LD05246PASB A/C

COMPOUND	SPIKE ADDED (ug/kg)	SAMPLE CONCENTRATION (ug/kg)	LCS CONCENTRATION (ug/kg)	LCS RECOVERY %	#	QC LIMITS % REC
PCB 1016	267	U	244	91		67-144
PCB 1260	267	U	224	84		58-155
PCB 1016 #2	267	U	268	100		69-140
PCB 1260 #2	267	U	279	104		67-149

COMPOUND	SPIKE ADDED (ug/kg)	LCSD CONCENTRATION (ug/kg)	LCSD % REC	#	RPD	#	QC LIMITS RPD	QC LIMITS % REC.
PCB 1016	267	255	95		4		30	67-144
PCB 1260	267	238	89		6		30	58-155
PCB 1016 #2	267	273	102		2		30	69-140
PCB 1260 #2	267	283	106		2		30	67-149

Column to be used to flag recovery and RPD values with an asterisk

* Values outside of QC limits
RPD: 0 out of 4 outside limits
Spike Recovery: 0 out of 8 outside limits

Comments: _____

analytics environmental laboratory LLC

PCB WIPE
LABORATORY CONTROL SPIKE/ DUPLICATE
PERCENT RECOVERY

Project #: 218736 Quote #: SDG#: 56617

Site ID: Location: ME

Instrument ID: F

GC Column: RTX-CLPesticides I ID: 0.32(mm)

RTX-CLPesticides II

Laboratory Control Spike/Duplicate Sample ID: L05256PWIPE/ LD05256PWIPE A/C

COMPOUND	SPIKE ADDED (ug/wipe)	SAMPLE CONCENTRATION (ug/wipe)	LCS CONCENTRATION (ug/wipe)	LCS RECOVERY % #	QC LIMITS % REC
PCB 1016	0.40	U	0.42	104	67-144
PCB 1260	0.40	U	0.41	101	58-155
PCB 1016 #2	0.40	U	0.44	109	69-140
PCB 1260 #2	0.40	U	0.45	113	67-149

COMPOUND	SPIKE ADDED (ug/wipe)	LCSD CONCENTRATION (ug/wipe)	LCSD % REC #	RPD #	QC LIMITS RPD	QC LIMITS % REC.
PCB 1016	0.40	0.41	102	2	30	67-144
PCB 1260	0.40	0.40	101	1	30	58-155
PCB 1016 #2	0.40	0.43	108	1	30	69-140
PCB 1260 #2	0.40	0.45	111	1	30	67-149

Column to be used to flag recovery and RPD values with an asterisk

* Values outside of QC limits
RPD: 0 out of 4 outside limits
Spike Recovery: 0 out of 8 outside limits

Comments: _____

PCB FORM 3

56617PCBF3.xls

analytics environmental laboratory LLC

PCB SOLID
LABORATORY CONTROL SPIKE
PERCENT RECOVERY

Project #: 218736 Quote #: SDG#: 56617

Site ID: Location: ME

Instrument ID: F

GC Column: RTX-CLPesticides I ID: 0.32(mm)

RTX-CLPesticides II

Laboratory Control Spike/Duplicate Sample ID: L05256PASB A/C

COMPOUND	SPIKE ADDED (ug/kg)	SAMPLE CONCENTRATION (ug/kg)	LCS CONCENTRATION (ug/kg)	LCS RECOVERY % #	QC LIMITS % REC
PCB 1016	267	U	272	102	67-144
PCB 1260	267	U	275	103	58-155
PCB 1016 #2	267	U	279	105	69-140
PCB 1260 #2	267	U	290	109	67-149

Column to be used to flag recovery and RPD values with an asterisk

* Values outside of QC limits

Spike Recovery: 0 out of 4 outside limits

Comments: _____

analytics environmental laboratory LLC

PCB SOLID
LABORATORY CONTROL SPIKE/ DUPLICATE
PERCENT RECOVERY

Project #: 218736 Quote #: SDG#: 56617

Site ID: Location: ME

Instrument ID: F

GC Column: RTX-CLPesticides I ID: 0.32(mm)

RTX-CLPesticides II

Laboratory Control Spike/Duplicate Sample ID: L05266PASB A/C/ LD05266PASB A/C

COMPOUND	SPIKE ADDED (ug/kg)	SAMPLE CONCENTRATION (ug/kg)	LCS CONCENTRATION (ug/kg)	LCS RECOVERY %	#	QC LIMITS % REC
PCB 1016	267	U	260	97		67-144
PCB 1260	267	U	247	92		58-155
PCB 1016 #2	267	U	270	101		69-140
PCB 1260 #2	267	U	280	105		67-149

COMPOUND	SPIKE ADDED (ug/kg)	LCSD CONCENTRATION (ug/kg)	LCSD % REC	#	RPD	#	QC LIMITS RPD	QC LIMITS % REC.
PCB 1016	267	274	103		5		30	67-144
PCB 1260	267	269	101		9		30	58-155
PCB 1016 #2	267	285	107		5		30	69-140
PCB 1260 #2	267	296	111		6		30	67-149

Column to be used to flag recovery and RPD values with an asterisk

* Values outside of QC limits
RPD: 0 out of 4 outside limits
Spike Recovery: 0 out of 8 outside limits

Comments: _____

analytics environmental laboratory LLC

PCB SOLID
MATRIX SPIKE/ DUPLICATE
PERCENT RECOVERY

Project #: 218736 Quote #: SDG#: 56617

Site ID: Location: ME

Instrument ID: F

GC Column: RTX-CL Pesticides I ID: 0.32(mm)

RTX-CL Pesticides II

Matrix Spike/Duplicate Sample ID: 56617-23

COMPOUND	SPIKE ADDED (ug/kg)	SAMPLE CONCENTRATION (ug/kg)	MS CONCENTRATION (ug/kg)	MS RECOVERY %	#	QC LIMITS % REC
PCB 1016	297	U	281	95		67-144
PCB 1260	297	**49**	454	136		58-155
PCB 1016 #2	297	U	295	99		69-140
PCB 1260 #2	297	**50**	458	137		67-149

COMPOUND	SPIKE ADDED (ug/kg)	MSD CONCENTRATION (ug/kg)	MSD % REC	#	RPD	#	QC LIMITS RPD	QC LIMITS % REC.
PCB 1016	342	336	98		18		30	67-144
PCB 1260	342	362	92		22		30	58-155
PCB 1016 #2	342	353	103		18		30	69-140
PCB 1260 #2	342	366	92		22		30	67-149

Column to be used to flag recovery and RPD values with an asterisk

* Values outside of QC limits
RPD: 0 out of 4 outside limits
Spike Recovery: 0 out of 8 outside limits

Comments: _____

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PEST/PCB/HERB
Instrument F
Sequence Log Book

051606 -F

Load: [signature]
Unload: [signature]

Data File	Vial #	Sample	Method	Comments
F55857	9	P4045 Pest50		✓
58	10	P4046 Pest100		✓
59	11	P4047 Pest150		✓
60	12	P3965 Pest 0.5CC		✓
61	13	P4016 16/60 0.01	PBQ051166	✓
62	14	P4017 16/60 0.02		✓
63	15	P4018 16/60 0.05		✓
64	16	P4019 16/60 0.1		✓
65	17	P4020 16/60 0.25		✓
66	18	P4021 16/60 0.5		✓
67	19	P4022 16/60 1.0		✓
68	20	P4024 16/60 0.5CC		✓
69	21	P3883 AR1221		for overlay, single point
70	22	P3884 AR1232		
71	23	P3885 AR 1242		
72	24	P3886 AR 1248		
73	25	P3994 AR 1254		
74	26	P3887 AR 1262		
75	27	P3888 AR 1268	5.18.06	
76	28	P4057 AR1254 0.02	P540051166	✓
77	29	P4056 AR1254 0.05		✓
78	30	P4055 AR1254 0.1		✓
79	31	P4054 AR1254 0.25		✓
80	32	P4053 AR1254 0.5		✓
81	33	P4052 AR1254 1.0		✓
82	34	P4010 AR1254 0.5		✓
83	35	B051106PW		✓
84	36	L051106PWT		✓ methoxy ↑#1, ok#2, endrin ketone ↑(marg.)#1, ↑#2
85	37	L051106PWB		✓ 5.17.06

Standard Dilutions: 1:2- 500 uL sample +500 uL hexane
1:5- 200 uL sample +800 uL hexane
Hexane Lot : _____ 1:10- 100 uL sample + 900 uL hexane
1:20- 50 uL sample + 950 uL hexane
MTBE Lot: _____ 1:100- 10 uLsample +990 uL hexane
1:1000- 1 uLsample +1000 uL hexane

PEST/PCB/HERB
Instrument F
Sequence Log Book

Load: yy
Unload: yy

052406-F

Data File	Vial #	Sample	Method	Comments
F560056	100	Hexane		✓
57	1	P4D20 16/160 0.25	PBQ05166	✓
58	2	B052464PAS A/C		✓
59	3	L052464PASB		✓
60	4	LD052464PASB		✓
61	5	56606-1		✓
62	6	-2		✓
63	7	-3		✓
64	8	-4		✓
65	9	-5		RR@ 1:10
66	10	-6		RR@ 1:2000
67	11	-6.MS		↑1260 hit in sample
68	12	-6.MSD		ms/msD failed
69	13	P4021 16/160 0.5		✓

5.25.00

Standard Dilutions: 1:2- 500 uL sample +500 uL hexane
1:5- 200 uL sample +800 uL hexane
Hexane Lot : _____ 1:10- 100 uL sample + 900 uL hexane
1:20- 50 uL sample + 950 uL hexane
MTBE Lot: _____ 1:100- 10 uLsample +990 uL hexane
1:1000- 1 uLsample +1000 uL hexane

Analytics environmental laboratory LLC

PEST/PCB/HERB
Instrument F
Sequence Log Book

052506-F

Load: YJ
Unload: YJ

Data File	Vial #	Sample	Method	Comments
F56070	1	P4020 16|60 0.25		not diluting would peak.
71	100	Hexane		sys. conditioning
72	1	P4020 16|60 0.25	PBQ051666	✓
73	2	B052466PAS A|C		
74	14	566006-5, 1:10		✓
75	15	566006-6, 1:2000 ↓		✓
76	13	P4021 16|60 0.5		✓
77	100	Hexane		changed liner
~~78~~	~~100~~	~~Hex~~		YJ 5.25.06
78	1	P4020 16|60 0.25	PBQ051666	✓
79	2	B052566P Wipe, A|C		✓
80	3	L052566P Wipe		✓
81	4	LD052566P Wipe ↓		✓
82	5	56617-24		✓
83	6	-25		✓
84	7	-26		✓
85	8	-27		✓
86	9	-28		RR@1:10
87	10	-29		✓
88	11	-30		✓
89	12	-31		✓
90	13	-32		✓
91	14	-33		✓
92	15	-34		✓
93	16	-35		✓
94	17	-36		✓
95	18	-37 ↓		✓
96	19	P4021 16|60 0.5		✓
97	20	B052566PAS A|C		✓

Standard Dilutions: 1:2- 500 uL sample +500 uL hexane
1:5- 200 uL sample +800 uL hexane
Hexane Lot : _____ 1:10- 100 uL sample + 900 uL hexane
1:20- 50 uL sample + 950 uL hexane
MTBE Lot: _____ 1:100- 10 uLsample +990 uL hexane
1:1000- 1 uLsample +1000 uL hexane

analytics environmental laboratory LLC

PEST/PCB/HERB
Instrument F
Sequence Log Book

052506-F

Load: JJ
Unload: JJ

Data File	Vial #	Sample	Method	Comments
F56098	21	L052506PASB A|C	PB1205166	✓
99	22	56617-1		✓
F56100	23	-2	54505166	✓ 1260/1254
01	24	-6		✓
02	25	-7		✓
03	26	-8		✓
04	27	-9		✓
05	28	-10		✓
06	29	-11		✓
07	30	-12		✓
08	31	-13	42505166	✓ 1242
~~09~~	~~32~~			JJ 5.25.06
09	19	P4021 16160 0.5		✓
10	32	56617-14 A|C		RR@1:2000
11	33	56617-3, 1:10		RR@1:5
12	34	-4, 1:10	42505166	✓ 1242
13	35	-5		✓
~~JJ 5.25.06~~ 14	19	P4021 16160 0.5		✓
15	100	Hexane		syst cond.
16	36	BO52506PAS A|C		✓
17	37	LO52506PASB		✓
18	38	LDO52506PASB		✓
19	39	56617-15		✓
20	40	-16		RR@1:10 1260/1254
21	41	-17		1260 RR@1:50 1254 RR@1:50
22	42	-18		✓
23	43	-19		✓
24	44	-20		✓
25	45	-21		✓

Standard Dilutions: 1:2- 500 uL sample +500 uL hexane
1:5- 200 uL sample +800 uL hexane
Hexane Lot : 46060 1:10- 100 uL sample + 900 uL hexane
1:20- 50 uL sample + 950 uL hexane
MTBE Lot: _____ 1:100- 10 uLsample +990 uL hexane
1:1000- 1 uLsample +1000 uL hexane

~~1:50 - 20ul sample 980 ul hexane~~ JJ 5.26.06

1:2000 = 1ul
2000ul

analytics environmental laboratory LLC

PEST/PCB/HERB
Instrument F
Sequence Log Book

052506-F

Load: JJ 5.26.06
Unload: JJ

Data File	Vial #	Sample	Method	Comments
F56126	46	56617-22	PBQD5166	✓
27	47	-23		✓
28	48	-23, MS		✓ used actual chrts for
29	49	-23, MSD		✓ RPD values, OK.
30	19	P4021 16/60 D5		✓

Standard Dilutions: 1:2- 500 uL sample +500 uL hexane
1:5- 200 uL sample +800 uL hexane
Hexane Lot : _____ 1:10- 100 uL sample + 900 uL hexane
1:20- 50 uL sample + 950 uL hexane
MTBE Lot: _____ 1:100- 10 uLsample +990 uL hexane
1:1000- 1 uLsample +1000 uL hexane

analytics environmental laboratory LLC

PEST/PCB/HERB
Instrument F
Sequence Log Book

052606-F

Load: JJ 5.26.06
Unload: JJ 5.26.06

Data File	Vial #	Sample	Method	Comments
F56131	1	P4020 16/60 0.25	PBQ051660	↑, susp. conc. repoured
32	50	P4020 16/600.25		✓ ↑ 1260 II 1254.
33 JJ 5.26.06	~~51~~ 36	~~56617-28~~, BQ5266PPSA/C JJ 5.26.06		✓
34	51	56617-28, 1:10 A/C		✓
35	52	-14, 1:2000		✓
36	53	-3, 1:5	545051660	✓ 1260/1254
37	54	-16, 1:10		✓ 1260/1254
38	55	-17, 1:50		✓ 1260/1254
39	19	P4021 16/60 0.5		✓

* window continued on next page. I'd out before complete

JJ 5.26.06

Standard Dilutions: 1:2- 500 uL sample +500 uL hexane
1:5- 200 uL sample +800 uL hexane
Hexane Lot: 460120 1:10- 100 uL sample + 900 uL hexane
1:20- 50 uL sample + 950 uL hexane
MTBE Lot: _____ 1:100- 10 uLsample +990 uL hexane
1:1000- 1 uLsample +1000 uL hexane

1:2000 - 1ul into 2.0ml
1:50 - 20ul into 980ul

C:\HPCHEM\2\METHODS\PBQ05166.M

Level 1 File Name	F55861.D
Level 2 File Name	F55862.D
Level 3 File Name	F55863.D
Level 4 File Name	F55864.D
Level 5 File Name	F55865.D
Level 6 File Name	F55866.D
Level 7 File Name	F55867.D

5-17-06

5.17.06

RESPONSE = X + YA + Z(A*A)

Name	R^2	R	X	Y	Z
TCX	1.000	1.000	1.04E+04	3.47E+07	0.00E-01
Ar1016 A	0.999	0.999	2.45E+04	1.65E+06	0.00E-01
Ar1016 B	0.999	1.000	3.80E+04	3.18E+06	0.00E-01
Ar1016 C	1.000	1.000	1.15E+04	1.34E+06	0.00E-01
Ar1260 A	0.999	1.000	2.75E+04	2.04E+06	0.00E-01
Ar1260 B	1.000	1.000	2.73E+04	2.45E+06	0.00E-01
Ar1260 C	1.000	1.000	8.21E+03	1.96E+06	0.00E-01
Ar1260 D	1.000	1.000	7.55E+03	1.93E+06	0.00E-01
DCB	0.999	0.999	9.77E+04	2.91E+07	0.00E-01
TCX #2	0.999	1.000	6.78E+04	3.70E+07	0.00E-01
Ar1016 A #2	0.997	0.999	4.34E+04	1.69E+06	0.00E-01
Ar1016 B #2	0.998	0.999	8.89E+04	3.83E+06	0.00E-01
Ar1016 C #2	0.997	0.999	3.37E+04	1.45E+06	0.00E-01
Ar1260 A #2	0.995	0.998	8.05E+04	2.28E+06	0.00E-01
Ar1260 B #2	0.996	0.998	1.10E+05	3.42E+06	0.00E-01
Ar1260 C #2	0.997	0.999	6.09E+04	2.54E+06	0.00E-01
Ar1260 D #2	0.999	0.999	7.76E+04	5.97E+06	0.00E-01
DCB #2	0.994	0.997	2.64E+05	3.87E+07	0.00E-01

Quantitation Report (QT Reviewed)

Signal #1 : C:\HPCHEM\2\DATA\051606-F\F55861.D\ECD1A.CH Vial: 13
Signal #2 : C:\HPCHEM\2\DATA\051606-F\F55861.D\ECD2B.CH
Acq On : 16 May 2006 7:38 pm Operator:
Sample : P4016 16/60 0.01 Inst : Instrumen
Misc : Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 17 9:56 2006 Quant Results File: PBQ05166.RES

Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 09:54:21 2006
Response via : Initial Calibration
DataAcq Meth : PSTQUICK.M

5.17.06 5.17.06

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film

Compound	RT#1	RT#2	Resp#1	Resp#2	ug/ml	ug/ml
System Monitoring Compounds						
1) S TCX	2.64	2.72	61307	77106	0.002	0.001 #
Spiked Amount 0.100	Range 46 - 122		Recovery =		2.00%#	1.00%#
9) S DCB	10.39	10.04	79033	133576	N.D. m	N.D. m
Spiked Amount 0.100	Range 26 - 86		Recovery =		0.00%#	0.00%#
Target Compounds						
2) L1 Ar1016 A	3.43	3.37	19531	27177	N.D. m	N.D.
3) L1 Ar1016 B	3.92	3.86	43320	57040	0.001	N.D. #
4) L1 Ar1016 C	4.09	4.03	15699	21986	0.004	N.D. #
Sum Ar1016 A			59019	0	0.001	N.D.
Average Ar1016 A					0.000	0.000
5) L2 Ar1260 A	6.67	6.35	30275	39717	0.002	N.D. m#
6) L2 Ar1260 B	7.03	6.81	33089	60122	0.001	N.D. m#
7) L2 Ar1260 C	7.92	7.53	22273	35850	0.006	N.D. m#
8) L2 Ar1260 D	8.49	8.59	22028	68293	0.004	N.D. #
Sum Ar1260 A			107665	0	0.013	N.D.
Average Ar1260 A					0.003	0.000

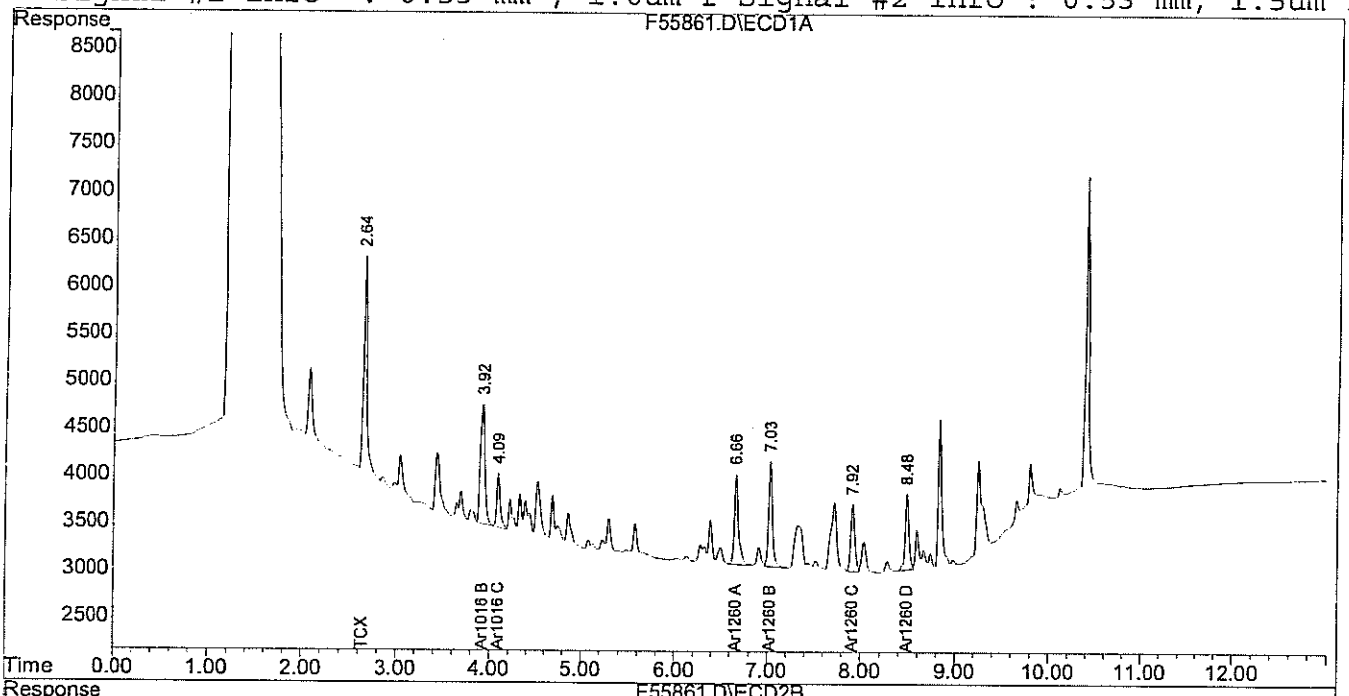
(f)=RT Delta > 1/2 Window (#)=Amounts differ by > 25% (m)=manual int

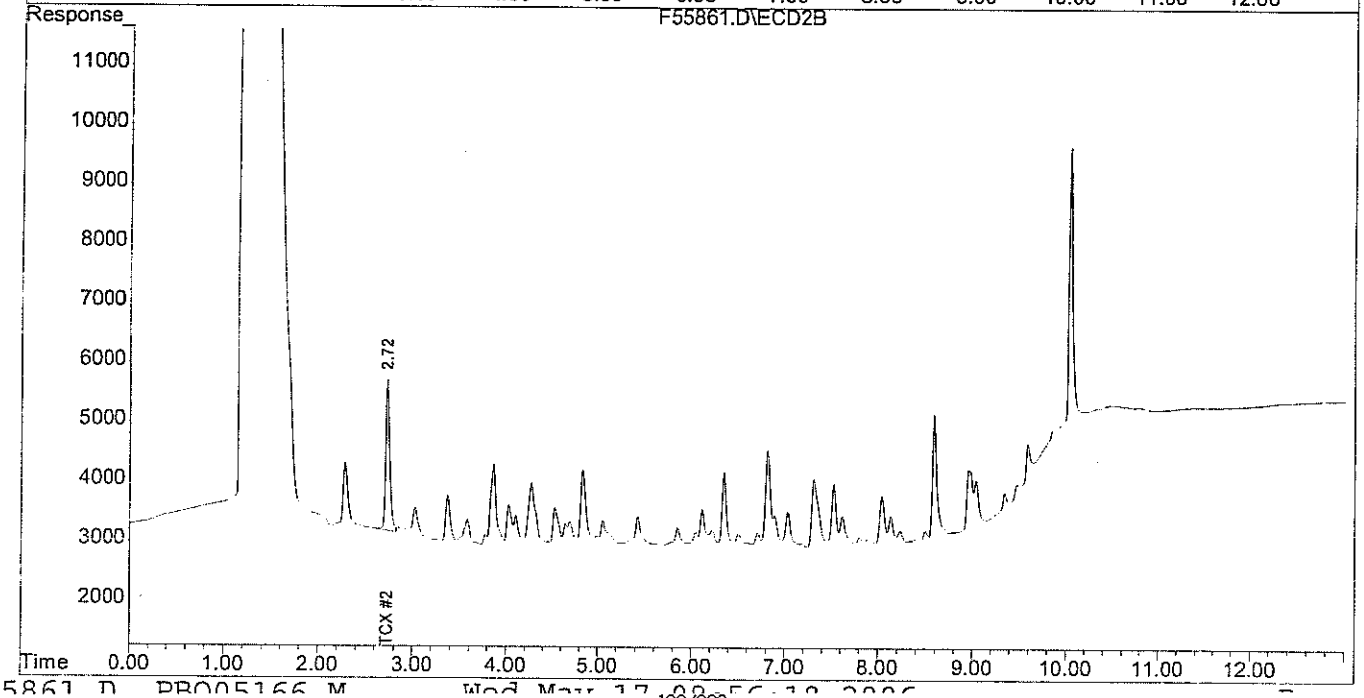
Quantitation Report (QT Reviewed)

Signal #1 : C:\HPCHEM\2\DATA\051606-F\F55861.D\ECD1A.CH Vial: 13
Signal #2 : C:\HPCHEM\2\DATA\051606-F\F55861.D\ECD2B.CH
Acq On : 16 May 2006 7:38 pm Operator:
Sample : P4016 16/60 0.01 Inst : Instrumen
Misc : Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 17 9:56 2006 Quant Results File: PBQ05166.RES

Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 09:54:21 2006
Response via : Multiple Level Calibration
DataAcq Meth : PSTQUICK.M

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film



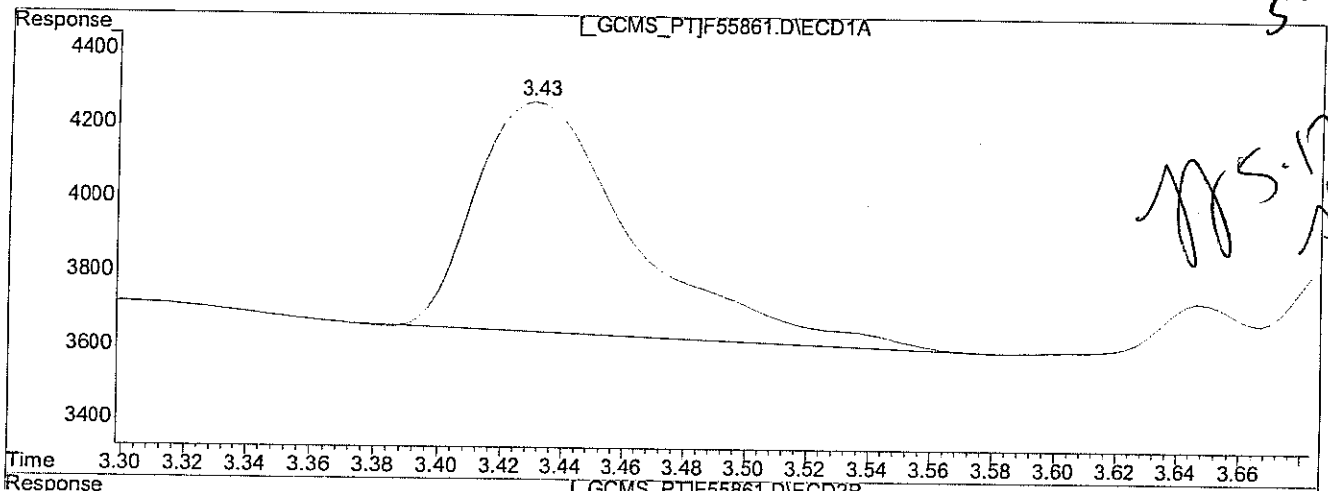


Quantitation Report (Qedit)

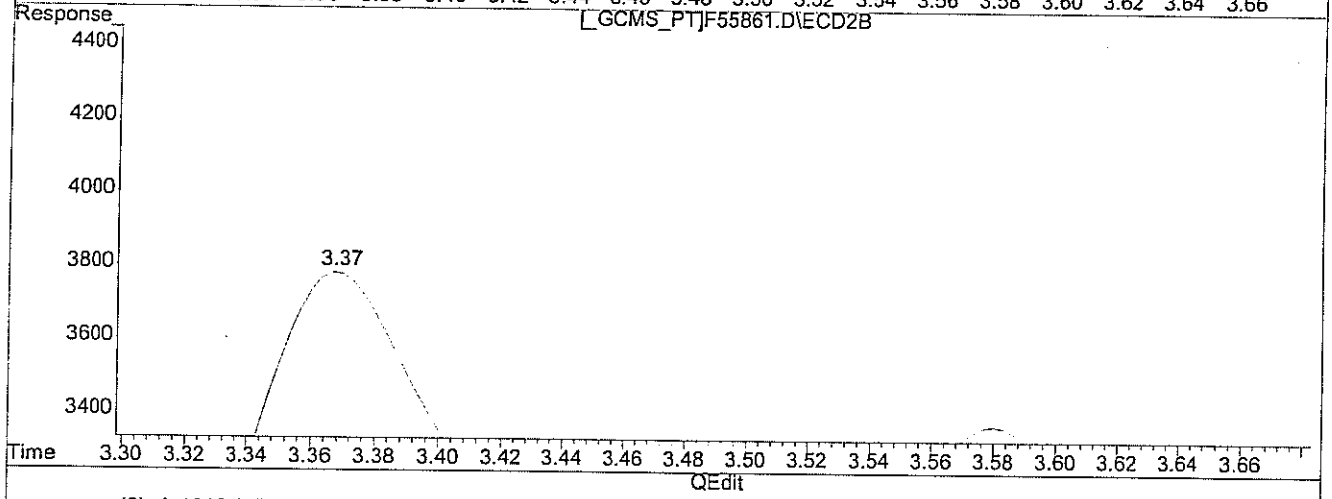
Signal #1 : C:\HPCHEM\2\DATA\051606-F\F55861.D\ECD1A.CH Vial: 13
Signal #2 : C:\HPCHEM\2\DATA\051606-F\F55861.D\ECD2B.CH
Acq On : 16 May 2006 7:38 pm Operator:
Sample : P4016 16/60 0.01 Inst : Instrumen
Misc : Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 17 9:54 2006 Quant Results File: PBQ05166.RES

Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 09:54:21 2006
Response via : Multiple Level Calibration

5.17.06



5.17.06 Shoulder



(2) Ar1016 A (L1)

R.T.	Response	Conc
3.43	22374	-0.00
3.92	43320	0.00
4.09	15699	0.00

(2) Ar1016 A #2 (L1)

R.T.	Response	Conc
3.37	27177	-0.01
3.86	57040	-0.01
4.03	21986	-0.01

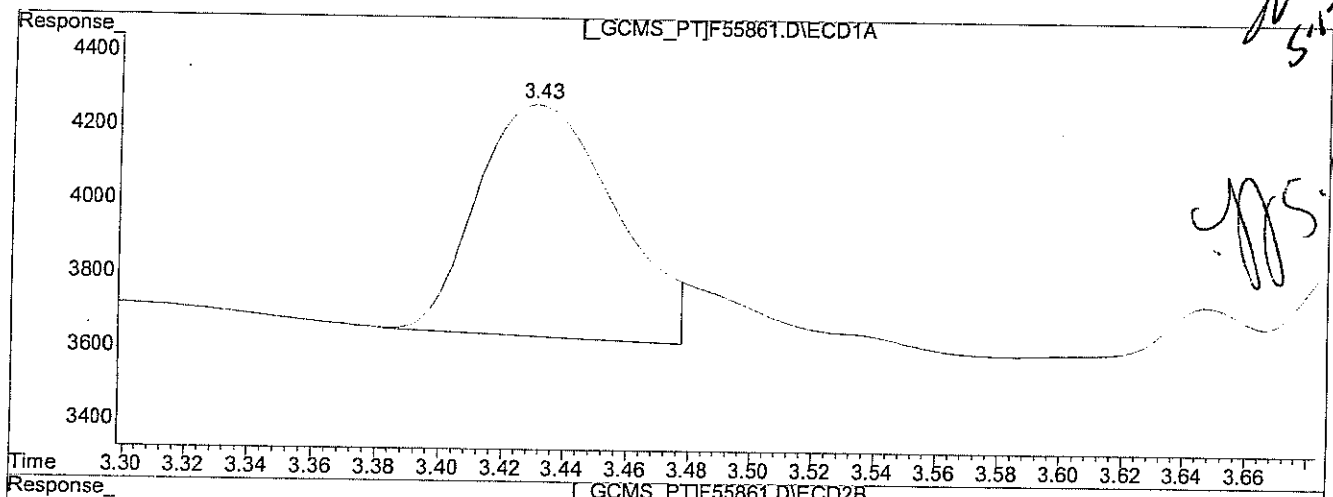
(+) = Expected Retention Time

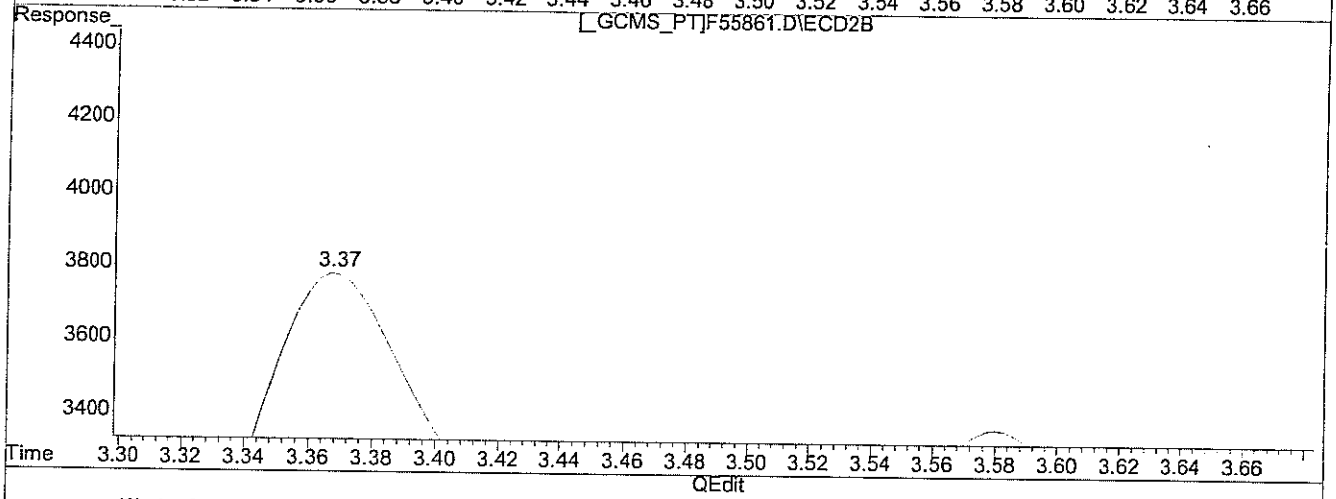
Quantitation Report (Qedit)

Signal #1 : C:\HPCHEM\2\DATA\051606-F\F55861.D\ECD1A.CH Vial: 13
Signal #2 : C:\HPCHEM\2\DATA\051606-F\F55861.D\ECD2B.CH
Acq On : 16 May 2006 7:38 pm Operator:
Sample : P4016 16/60 0.01 Inst : Instrumen
Misc : Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 17 9:54 2006 Quant Results File: PBQ05166.RES

Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 09:54:21 2006
Response via : Multiple Level Calibration

5.17.06





(2) Ar1016 A #2 (L1)

R.T.	Response	Conc
3.43	19531	-0.00
3.92	43320	0.00
4.09	15699	0.00

(2) Ar1016 A #2 (L1)

R.T.	Response	Conc
3.37	27177	-0.01
3.86	57040	-0.01
4.03	21986	-0.01

(+) = Expected Retention Time

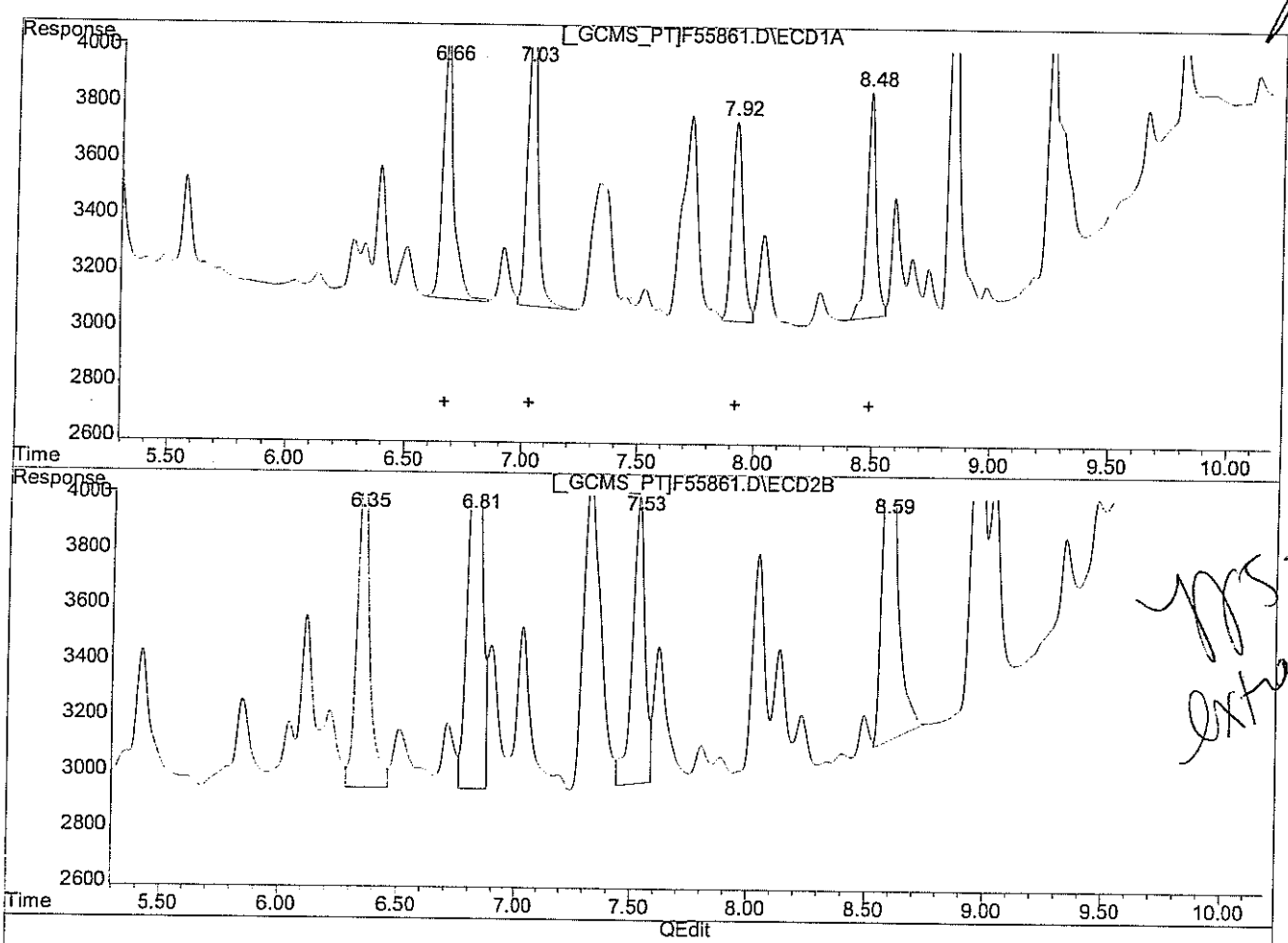
F55861.D PBQ05166.M Wed May 17 09:54:21 2006

Quantitation Report (Qedit)

Signal #1 : C:\HPCHEM\2\DATA\051606-F\F55861.D\ECD1A.CH Vial: 13
Signal #2 : C:\HPCHEM\2\DATA\051606-F\F55861.D\ECD2B.CH
Acq On : 16 May 2006 7:38 pm Operator:
Sample : P4016 16/60 0.01 Inst : Instrumen
Misc : Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 17 9:54 2006 Quant Results File: PBQ05166.RES

Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 09:54:21 2006
Response via : Multiple Level Calibration

5-17-06



5-17-06
extra area

(5) Ar1260 A (L2)

R.T.	Response	Conc
6.67	30275	0.00
7.03	33089	0.00
7.92	22273	0.01
8.49	22028	0.00

(5) Ar1260 A #2 (L2)

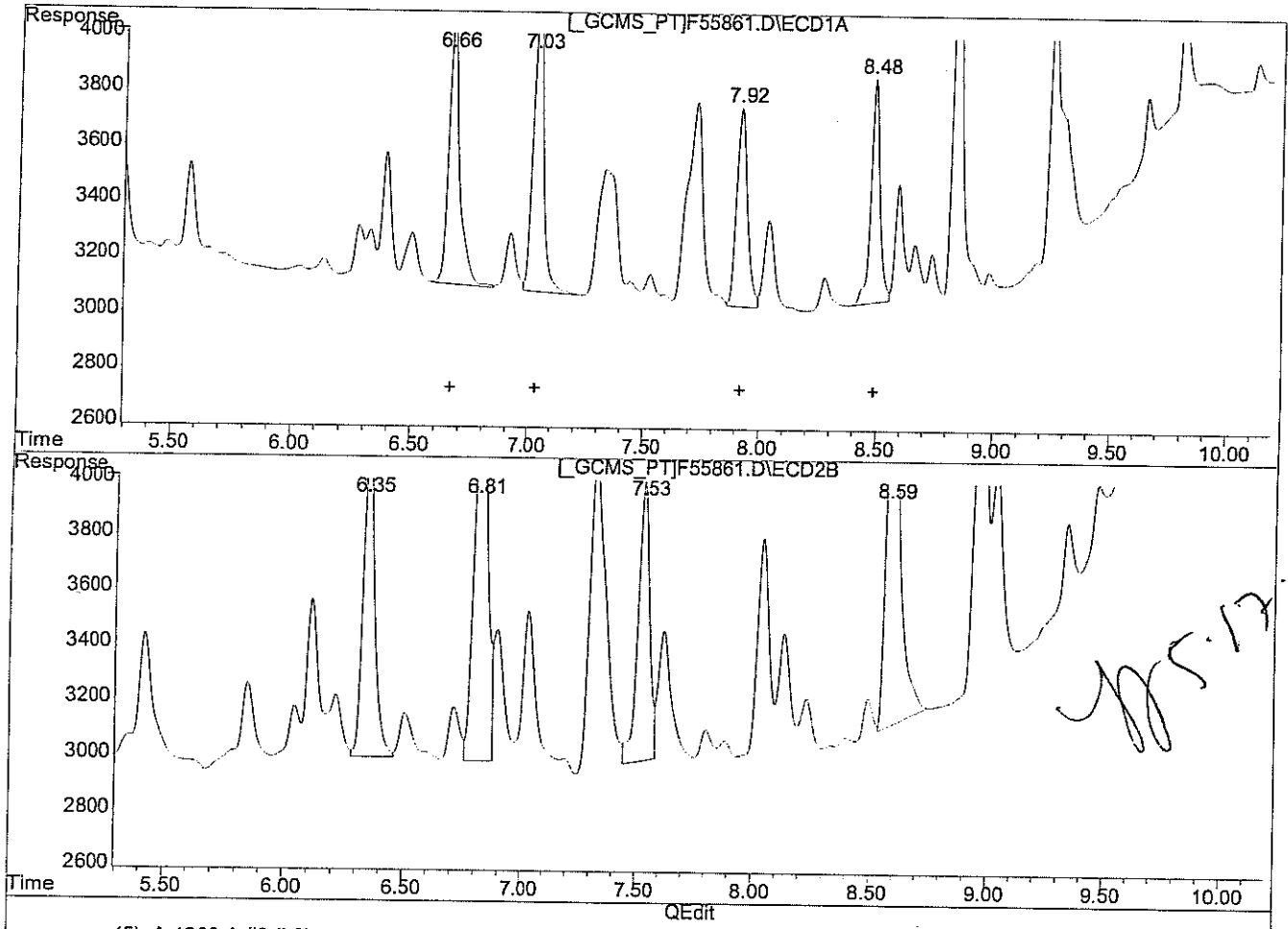
R.T.	Response	Conc
6.35	45036	-0.02
6.82	61803	-0.01
7.53	37673	-0.01
8.59	68293	-0.00

(+) = Expected Retention Time

Quantitation Report (Qedit)

Signal #1 : C:\HPCHEM\2\DATA\051606-F\F55861.D\ECD1A.CH Vial: 13
Signal #2 : C:\HPCHEM\2\DATA\051606-F\F55861.D\ECD2B.CH
Acq On : 16 May 2006 7:38 pm Operator:
Sample : P4016 16/60 0.01 Inst : Instrumen
Misc : Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 17 9:54 2006 Quant Results File: PBQ05166.RES

Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 09:54:21 2006
Response via : Multiple Level Calibration



(5) Ar1260 A #2 (L2)

R.T.	Response	Conc
6.67	30275	0.00
7.03	33089	0.00
7.92	22273	0.01
8.49	22028	0.00

(5) Ar1260 A #2 (L2)

R.T.	Response	Conc
6.35	39717	-0.02
6.81	60122	-0.01
7.53	35850	-0.01
8.59	68293	-0.00

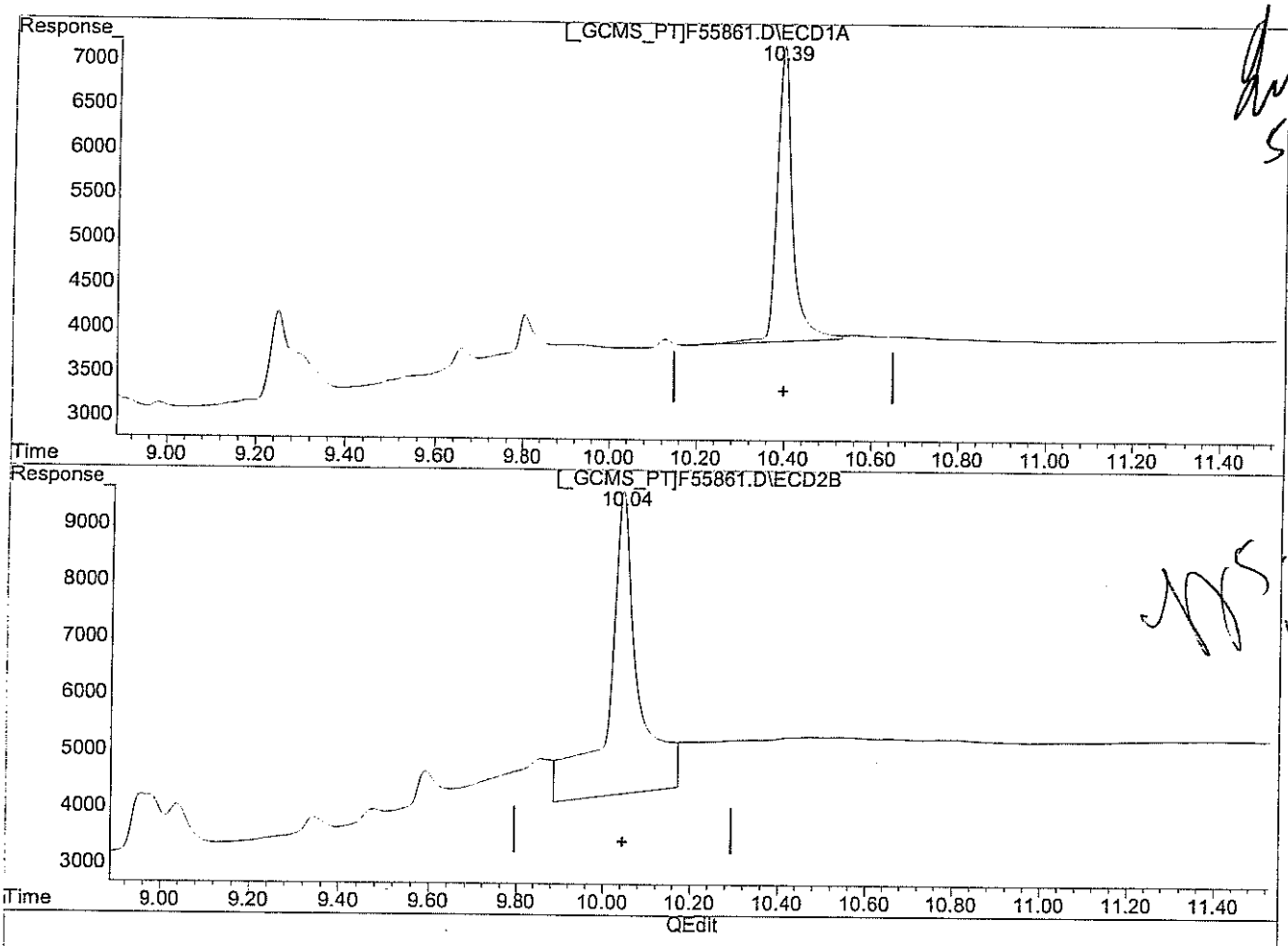
(+) = Expected Retention Time

F55861.D PBQ05166.M Wed May 17 09:55:50 2006

Quantitation Report (Qedit)

Signal #1 : C:\HPCHEM\2\DATA\051606-F\F55861.D\ECD1A.CH Vial: 13
Signal #2 : C:\HPCHEM\2\DATA\051606-F\F55861.D\ECD2B.CH
Acq On : 16 May 2006 7:38 pm Operator:
Sample : P4016 16/60 0.01 Inst : Instrumen
Misc : Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 17 9:54 2006 Quant Results File: PBQ05166.RES

Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 09:54:21 2006
Response via : Multiple Level Calibration



5.17.06

5.17.06 baseli

(9) DCB (S)

10.39min -0.001ug/ml

response 81717

(9) DCB #2 (S)

10.04min -0.000ug/ml

response 257756

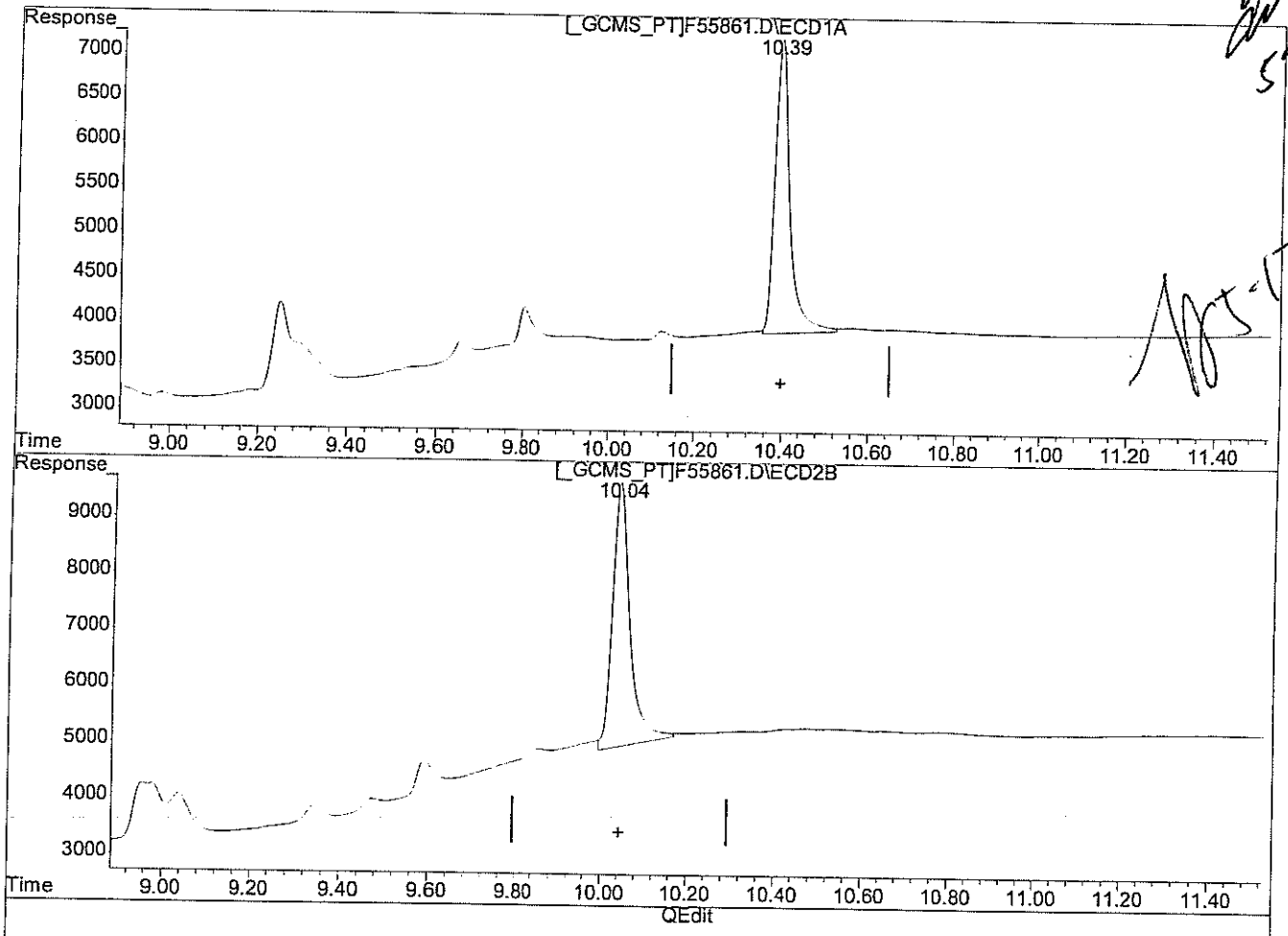
(+) = Expected Retention Time

F55861.D PBQ05166.M Wed May 17 09:55:02 2006

Quantitation Report (Qedit)

Signal #1 : C:\HPCHEM\2\DATA\051606-F\F55861.D\ECD1A.CH Vial: 13
Signal #2 : C:\HPCHEM\2\DATA\051606-F\F55861.D\ECD2B.CH
Acq On : 16 May 2006 7:38 pm Operator:
Sample : P4016 16/60 0.01 Inst : Instrumen
Misc : Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 17 9:54 2006 Quant Results File: PBQ05166.RES

Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 09:54:21 2006
Response via : Multiple Level Calibration



(9) DCB (S)

10.39min -0.001ug/ml m

response 79033

(9) DCB #2 (S)

10.04min -0.003ug/ml m

response 133576

(+) = Expected Retention Time

F55861.D PBQ05166.M

Quantitation Report (QT Reviewed)

Signal #1 : C:\HPCHEM\2\DATA\051606-F\F55862.D\ECD1A.CH Vial: 14
Signal #2 : C:\HPCHEM\2\DATA\051606-F\F55862.D\ECD2B.CH
Acq On : 16 May 2006 7:53 pm Operator:
Sample : P4017 16/60 0.02 Inst : Instrumen
Misc : Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 17 9:56 2006 Quant Results File: PBQ05166.RES

Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 09:54:21 2006
Response via : Initial Calibration
DataAcq Meth : PSTQUICK.M

5-17-06 5-17-06

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film

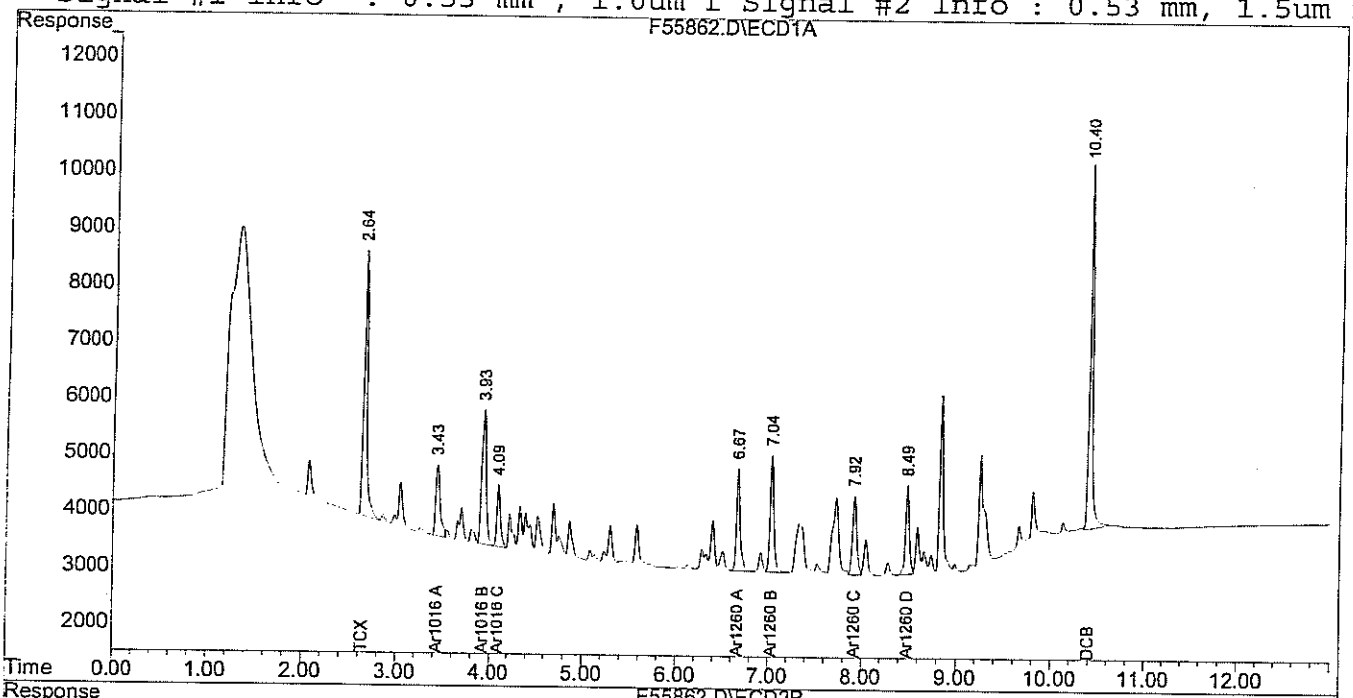
Compound	RT#1	RT#2	Resp#1	Resp#2	ug/ml	ug/ml
System Monitoring Compounds						
1) S TCX	2.64	2.73	122756	153548	0.004	0.003
Spiked Amount 0.100	Range 46 - 122			Recovery =	4.00%#	3.00%#
9) S DCB	10.40	10.04	155614	248581	0.001m	N.D. m#
Spiked Amount 0.100	Range 26 - 86			Recovery =	1.00%#	0.00%#
Target Compounds						
2) L1 Ar1016 A	3.44	3.37	42784	51825	0.011	0.006 #
3) L1 Ar1016 B	3.93	3.87	80946	111881	0.014	0.008 #
4) L1 Ar1016 C	4.09	4.03	29570	41472	0.016	0.008 #
Sum Ar1016 A			153300	205178	0.041	0.022
Average Ar1016 A					0.014	0.007
5) L2 Ar1260 A	6.67	6.35	53627	81296	0.014	0.002 #
6) L2 Ar1260 B	7.04	6.82	59571	113223	0.013	0.004 #
7) L2 Ar1260 C	7.92	7.54	42912	71915	0.017	0.007 #
8) L2 Ar1260 D	8.49	8.59	42372	139041	0.016	0.012 #
Sum Ar1260 A			198483	405475	0.060	0.024
Average Ar1260 A					0.015	0.006

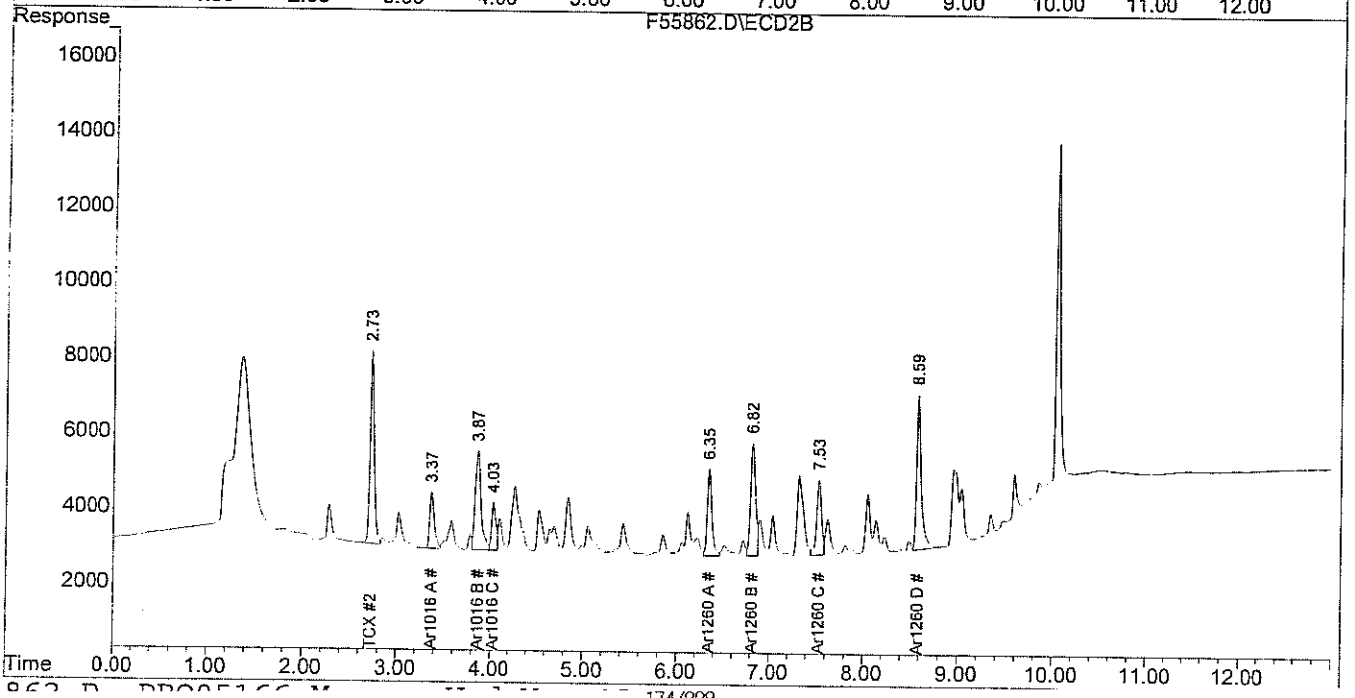
Quantitation Report (QT Reviewed)

Signal #1 : C:\HPCHEM\2\DATA\051606-F\F55862.D\ECD1A.CH Vial: 14
Signal #2 : C:\HPCHEM\2\DATA\051606-F\F55862.D\ECD2B.CH
Acq On : 16 May 2006 7:53 pm Operator:
Sample : P4017 16/60 0.02 Inst : Instrumen
Misc : Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 17 9:56 2006 Quant Results File: PBQ05166.RES

Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 09:54:21 2006
Response via : Multiple Level Calibration
DataAcq Meth : PSTQUICK.M

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film

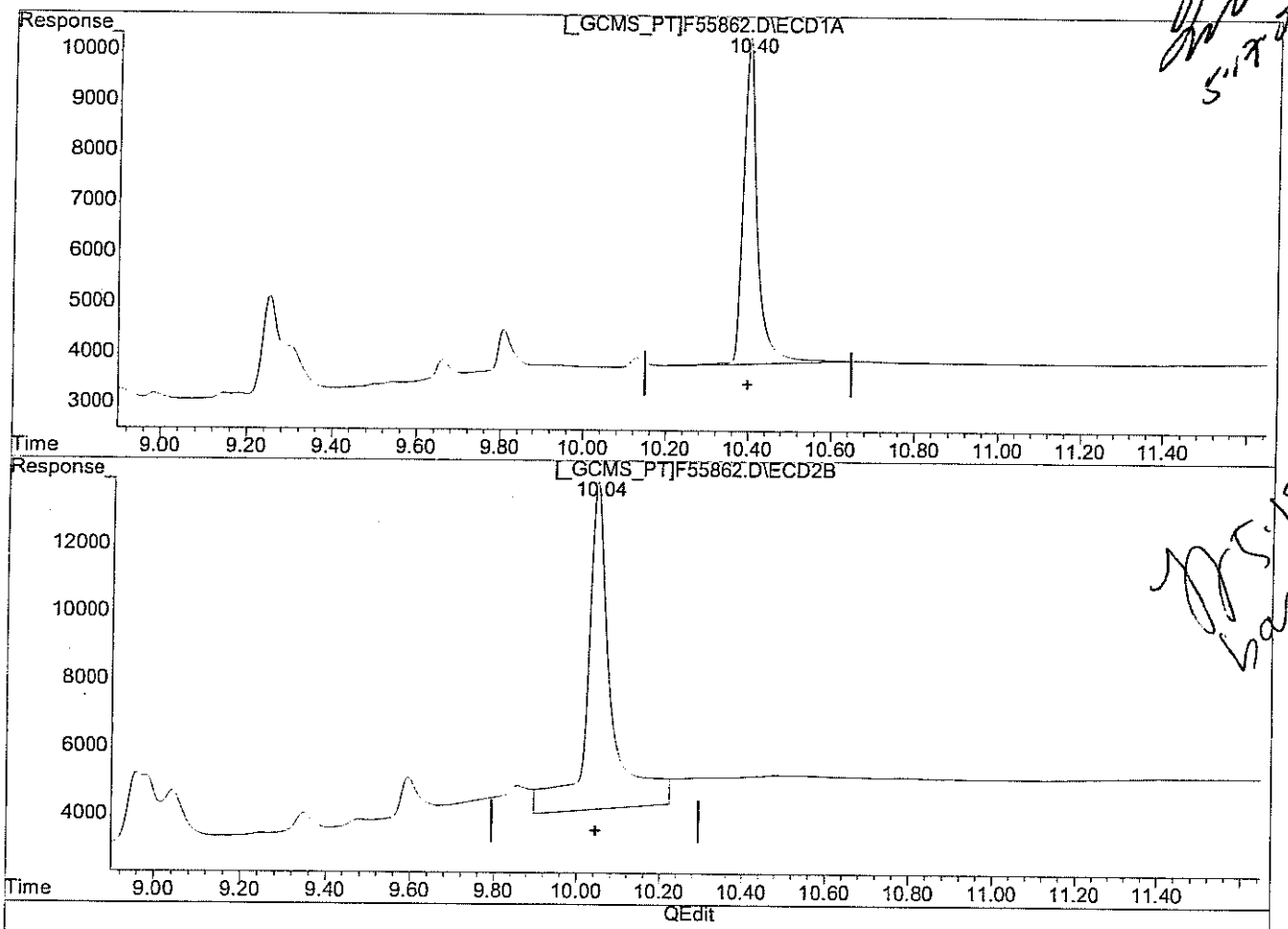




Quantitation Report (Qedit)

Signal #1 : C:\HPCHEM\2\DATA\051606-F\F55862.D\ECD1A.CH Vial: 14
Signal #2 : C:\HPCHEM\2\DATA\051606-F\F55862.D\ECD2B.CH
Acq On : 16 May 2006 7:53 pm Operator:
Sample : P4017 16/60 0.02 Inst : Instrumen
Misc : Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 17 9:56 2006 Quant Results File: PBQ05166.RES

Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 09:54:21 2006
Response via : Multiple Level Calibration



(9) DCB (S)

10.40min 0.002ug/ml

response 156938

(9) DCB #2 (S)

10.04min 0.003ug/ml

response 389832

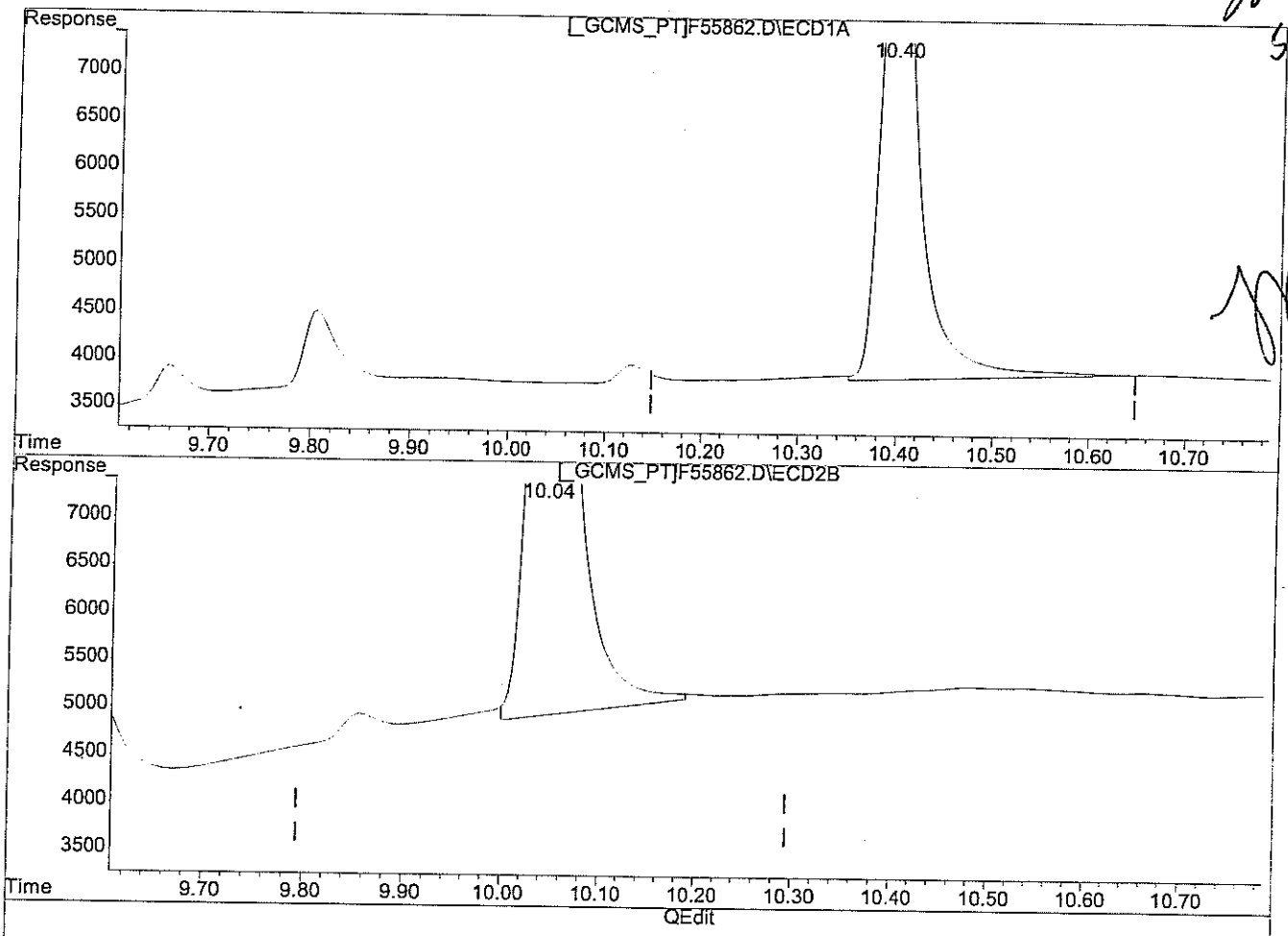
(+) = Expected Retention Time

Quantitation Report (Qedit)

Signal #1 : C:\HPCHEM\2\DATA\051606-F\F55862.D\ECD1A.CH Vial: 14
Signal #2 : C:\HPCHEM\2\DATA\051606-F\F55862.D\ECD2B.CH
Acq On : 16 May 2006 7:53 pm Operator:
Sample : P4017 16/60 0.02 Inst : Instrumen
Misc : Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 17 9:56 2006 Quant Results File: PBQ05166.RES

Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 09:54:21 2006
Response via : Multiple Level Calibration

5.17.06



5.17.06

(9) DCB (S)

10.40min 0.001ug/ml m

response 155614

(9) DCB #2 (S)

10.04min -0.000ug/ml m

response 248581

(+) = Expected Retention Time

F55862.D PBQ05166.M Wed M

Quantitation Report (QT Reviewed)

Signal #1 : C:\HPCHEM\2\DATA\051606-F\F55863.D\ECD1A.CH Vial: 15
Signal #2 : C:\HPCHEM\2\DATA\051606-F\F55863.D\ECD2B.CH
Acq On : 16 May 2006 8:07 pm Operator:
Sample : P4018 16/60 0.05 Inst : Instrumen
Misc : Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 17 10:04 2006 Quant Results File: PBQ05166.RES

Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 09:54:21 2006
Response via : Initial Calibration
DataAcq Meth : PSTQUICK.M

5.17.06 5.17.06

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film

Compound	RT#1	RT#2	Resp#1	Resp#2	ug/ml	ug/ml
System Monitoring Compounds						
1) S TCX	2.64	2.73	328762	401821	0.010	0.010
Spiked Amount 0.100	Range 46 - 122			Recovery =	10.00%#	10.00%#
9) S DCB	10.40	10.05	369688	589584	0.009m	0.009m
Spiked Amount 0.100	Range 26 - 86			Recovery =	9.00%#	9.00%#
Target Compounds						
2) L1 Ar1016 A	3.44	3.37	102888	121190	0.051	0.050
3) L1 Ar1016 B	3.93	3.87	187238	263561	0.051	0.050
4) L1 Ar1016 C	4.10	4.03	75426	99033	0.054	0.051
Sum Ar1016 A			365551	483784	0.156	0.151
Average Ar1016 A					0.052	0.050
5) L2 Ar1260 A	6.67	6.35	126226	190696	0.053	0.054
6) L2 Ar1260 B	7.04	6.82	148952	269972	0.053	0.054
7) L2 Ar1260 C	7.92	7.53	105396	183351	0.052	0.056
8) L2 Ar1260 D	8.49	8.59	102899	351344	0.051	0.051
Sum Ar1260 A			483473	995362	0.209	0.216
Average Ar1260 A					0.052	0.054

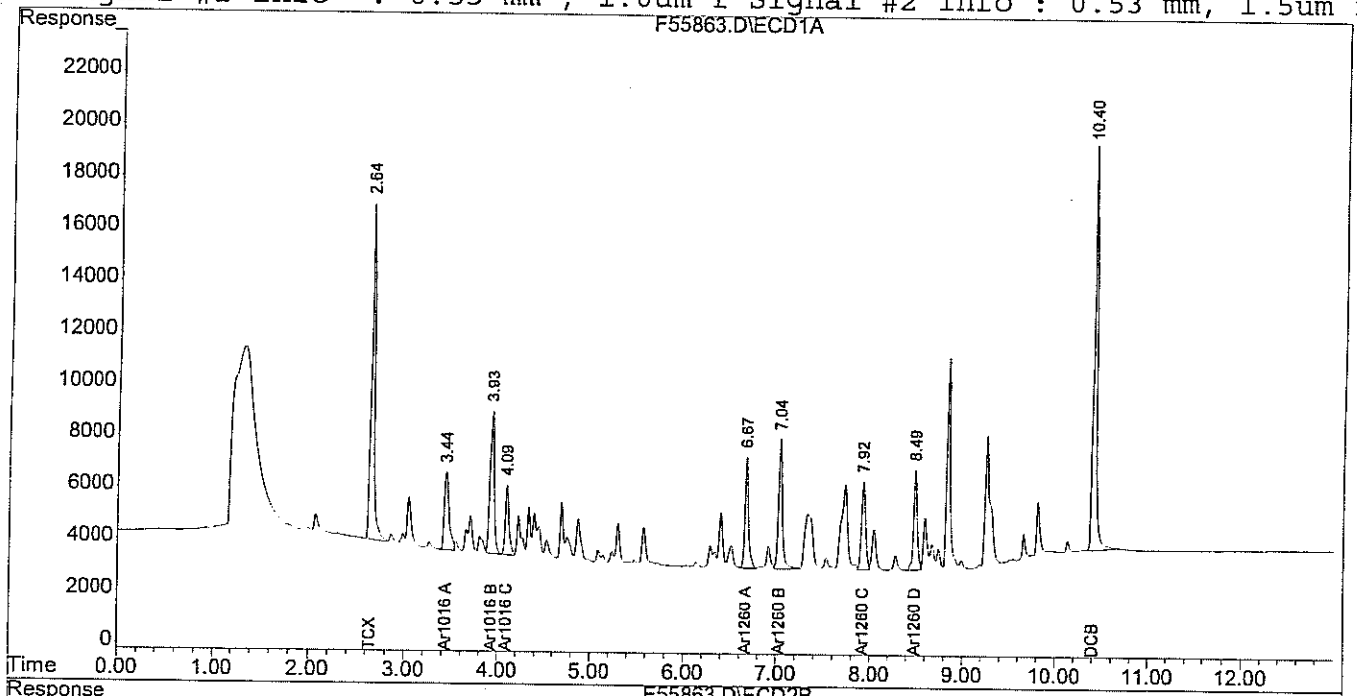
(f)=RT Delta > 1/2 Window (#)=Amounts differ by ... 25% / ...

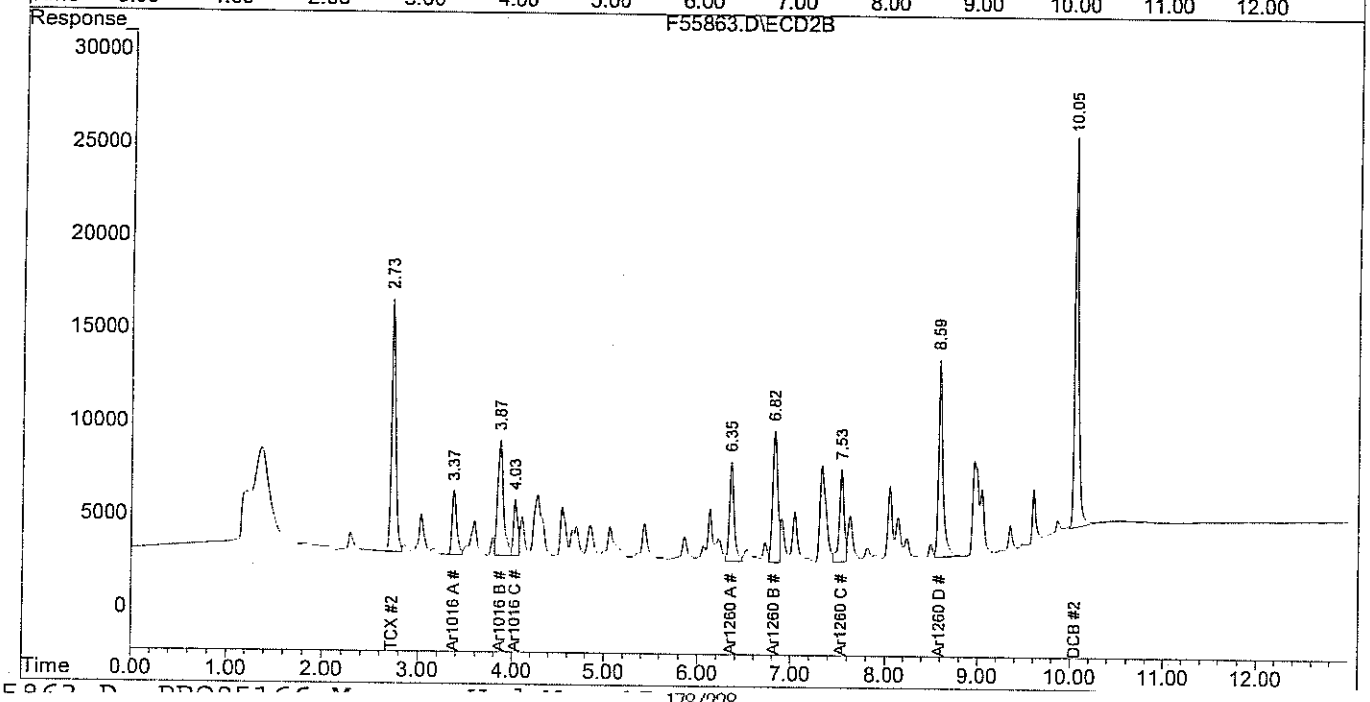
Quantitation Report (QT Reviewed)

Signal #1 : C:\HPCHEM\2\DATA\051606-F\F55863.D\ECD1A.CH Vial: 15
Signal #2 : C:\HPCHEM\2\DATA\051606-F\F55863.D\ECD2B.CH
Acq On : 16 May 2006 8:07 pm Operator:
Sample : P4018 16/60 0.05 Inst : Instrumen
Misc : Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 17 10:04 2006 Quant Results File: PBQ05166.RES

Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 09:54:21 2006
Response via : Multiple Level Calibration
DataAcq Meth : PSTQUICK.M

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film



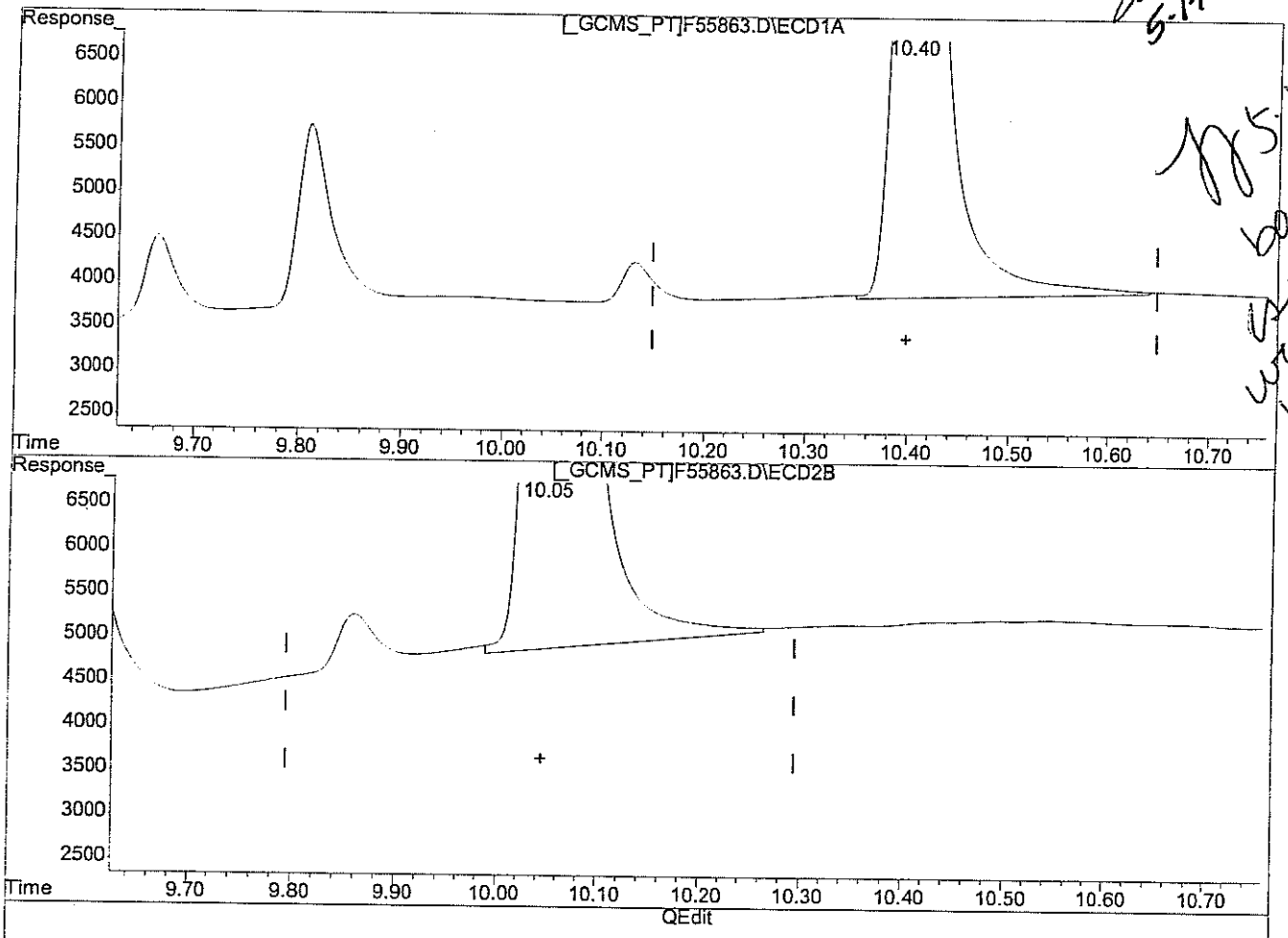


Quantitation Report (Qedit)

Signal #1 : C:\HPCHEM\2\DATA\051606-F\F55863.D\ECD1A.CH Vial: 15
Signal #2 : C:\HPCHEM\2\DATA\051606-F\F55863.D\ECD2B.CH
Acq On : 16 May 2006 8:07 pm Operator:
Sample : P4018 16/60 0.05 Inst : Instrumen
Misc : Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 17 9:57 2006 Quant Results File: PBQ05166.RES

Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 09:54:21 2006
Response via : Multiple Level Calibration

5-17-06



5-17-06 baseline issues were present before manual integration

(9) DCB (S)

10.40min 0.009ug/ml m

response 369688

(9) DCB #2 (S)

10.05min 0.009ug/ml m

response 589584

(+) = Expected Retention Time

F55863.D PBQ05166.M Wed May 17 09:54:21 2006

Quantitation Report (QT Reviewed)

Signal #1 : C:\HPCHEM\2\DATA\051606-F\F55864.D\ECD1A.CH Vial: 16
Signal #2 : C:\HPCHEM\2\DATA\051606-F\F55864.D\ECD2B.CH
Acq On : 16 May 2006 8:22 pm Operator:
Sample : P4019 16/60 0.1 Inst : Instrumen
Misc : Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 17 10:04 2006 Quant Results File: PBQ05166.RES

Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 09:54:21 2006
Response via : Initial Calibration
DataAcq Meth : PSTQUICK.M

5-17-06 5-17-06

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film

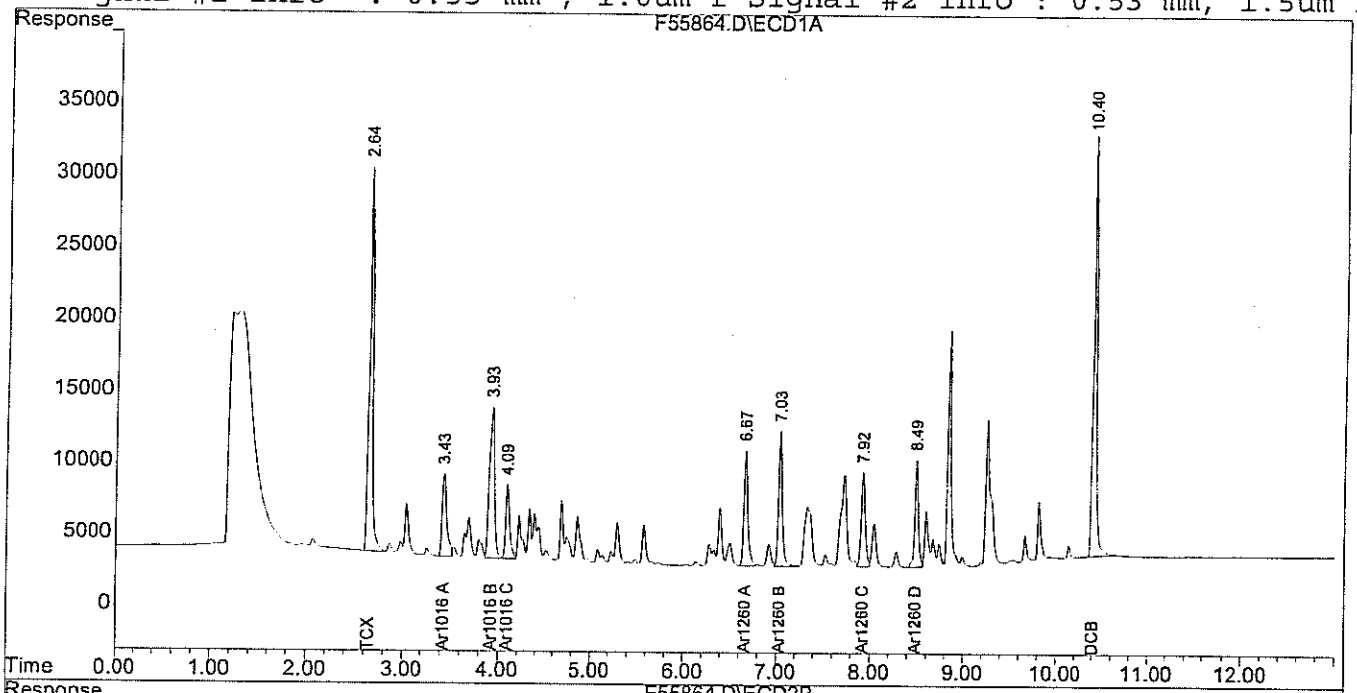
Compound	RT#1	RT#2	Resp#1	Resp#2	ug/ml	ug/ml
System Monitoring Compounds						
1) S TCX	2.64	2.73	688147	815855	0.022	0.022
Spiked Amount	0.100	Range 46 - 122		Recovery =	22.00%#	22.00%#
9) S DCB	10.40	10.05	698690	1099353	0.020	0.022m
Spiked Amount	0.100	Range 26 - 86		Recovery =	20.00%#	22.00%#
Target Compounds						
2) L1 Ar1016 A	3.43	3.37	199668	227033	0.116	0.117
3) L1 Ar1016 B	3.93	3.87	361008	502332	0.111	0.117
4) L1 Ar1016 C	4.09	4.03	146603	188678	0.113	0.118
Sum Ar1016 A			707279	918043	0.340	0.352
Average Ar1016 A					0.113	0.117
5) L2 Ar1260 A	6.67	6.35	234823	325259	0.112	0.119
6) L2 Ar1260 B	7.04	6.82	274639	476109	0.109	0.121
7) L2 Ar1260 C	7.92	7.53	203415	325317	0.107	0.119
8) L2 Ar1260 D	8.49	8.59	199734	691593	0.106	0.114
Sum Ar1260 A			912612	1818277	0.433	0.473
Average Ar1260 A					0.108	0.118

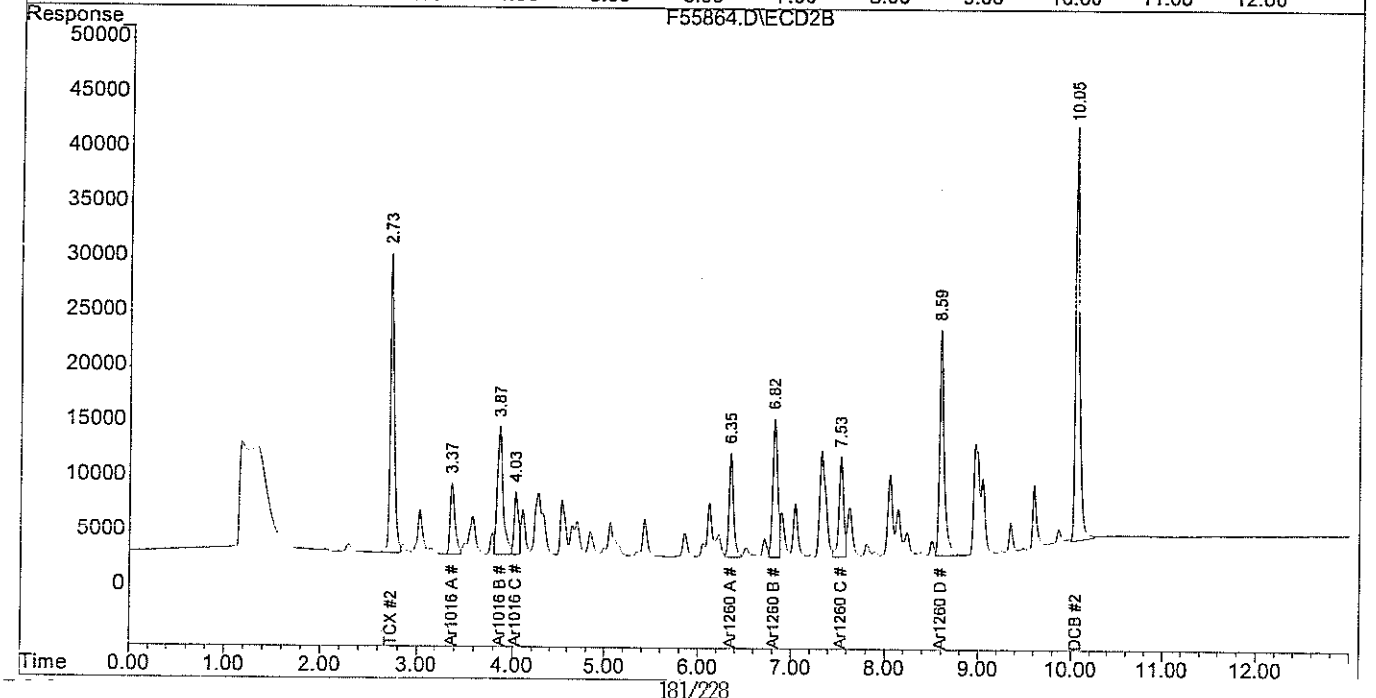
Quantitation Report (QT Reviewed)

Signal #1 : C:\HPCHEM\2\DATA\051606-F\F55864.D\ECD1A.CH Vial: 16
Signal #2 : C:\HPCHEM\2\DATA\051606-F\F55864.D\ECD2B.CH
Acq On : 16 May 2006 8:22 pm Operator:
Sample : P4019 16/60 0.1 Inst : Instrumen
Misc : Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 17 10:04 2006 Quant Results File: PBQ05166.RES

Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 09:54:21 2006
Response via : Multiple Level Calibration
DataAcq Meth : PSTQUICK.M

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film



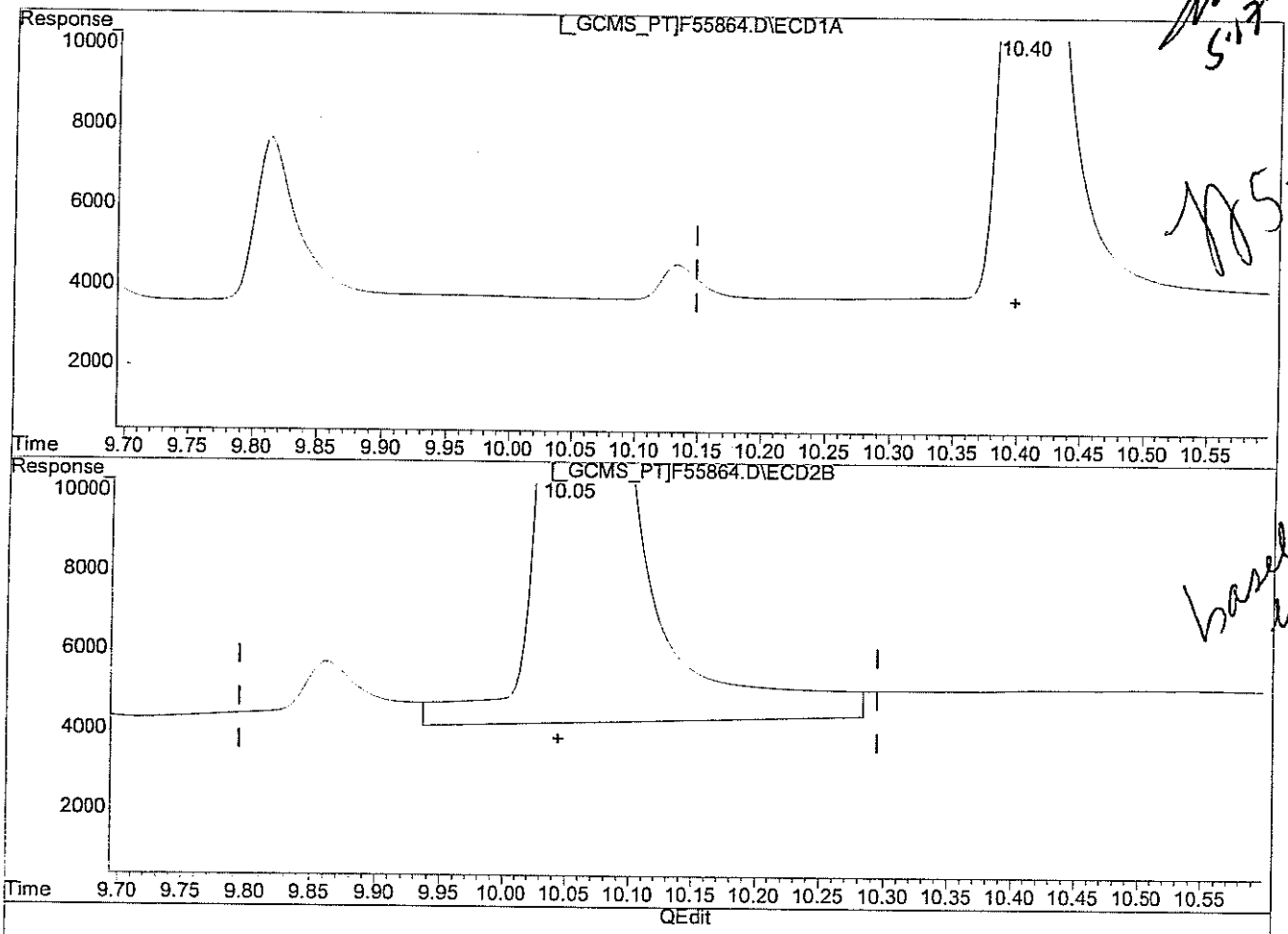


Quantitation Report (Qedit)

Signal #1 : C:\HPCHEM\2\DATA\051606-F\F55864.D\ECD1A.CH Vial: 16
Signal #2 : C:\HPCHEM\2\DATA\051606-F\F55864.D\ECD2B.CH
Acq On : 16 May 2006 8:22 pm Operator:
Sample : P4019 16/60 0.1 Inst : Instrumen
Misc : Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 17 10:04 2006 Quant Results File: PBQ05166.RES

Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 09:54:21 2006
Response via : Multiple Level Calibration

5-17-06



5-17-06

baseline/ extra area

(9) DCB (S)

10.40min 0.020ug/ml

response 698690

(9) DCB #2 (S)

10.05min 0.025ug/ml

response 1203022

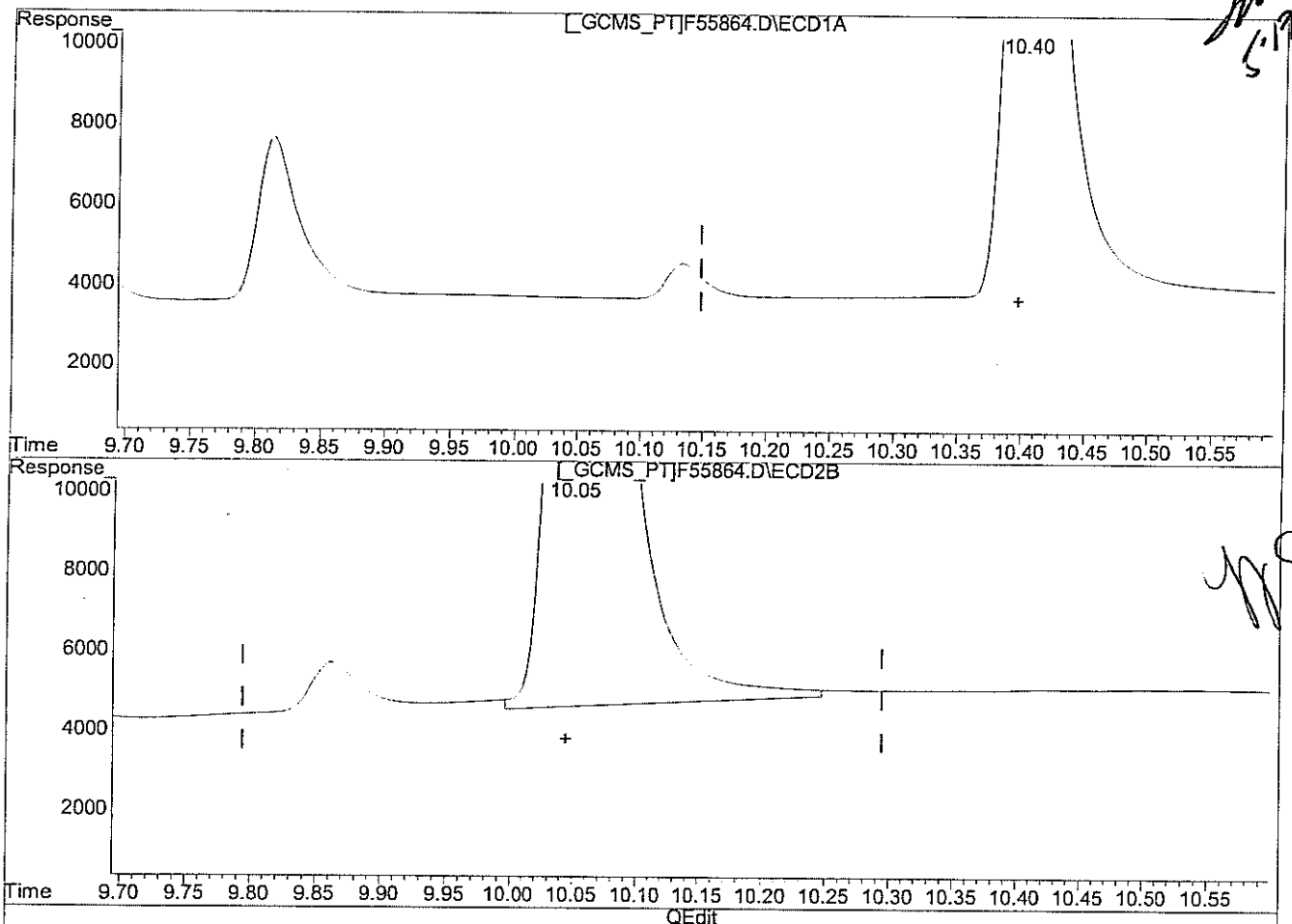
(+) = Expected Retention Time

Quantitation Report (Qedit)

Signal #1 : C:\HPCHEM\2\DATA\051606-F\F55864.D\ECD1A.CH Vial: 16
Signal #2 : C:\HPCHEM\2\DATA\051606-F\F55864.D\ECD2B.CH
Acq On : 16 May 2006 8:22 pm Operator:
Sample : P4019 16/60 0.1 Inst : Instrumen
Misc : Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 17 10:04 2006 Quant Results File: PBQ05166.RES

Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 09:54:21 2006
Response via : Multiple Level Calibration

5.17.06



5.17.06

(9) DCB (S)

10.40min 0.020ug/ml

response 698690

(9) DCB #2 (S)

10.05min 0.022ug/ml m

response 1099353

(+) = Expected Retention Time

Quantitation Report (QT Reviewed)

Signal #1 : C:\HPCHEM\2\DATA\051606-F\F55865.D\ECD1A.CH Vial: 17
Signal #2 : C:\HPCHEM\2\DATA\051606-F\F55865.D\ECD2B.CH
Acq On : 16 May 2006 8:36 pm Operator:
Sample : P4020 16/60 0.25 Inst : Instrumen
Misc : Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 17 9:54 2006 Quant Results File: PBQ05166.RES

Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 09:54:21 2006
Response via : Initial Calibration
DataAcq Meth : PSTQUICK.M

5.17.06 5.17.06

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film

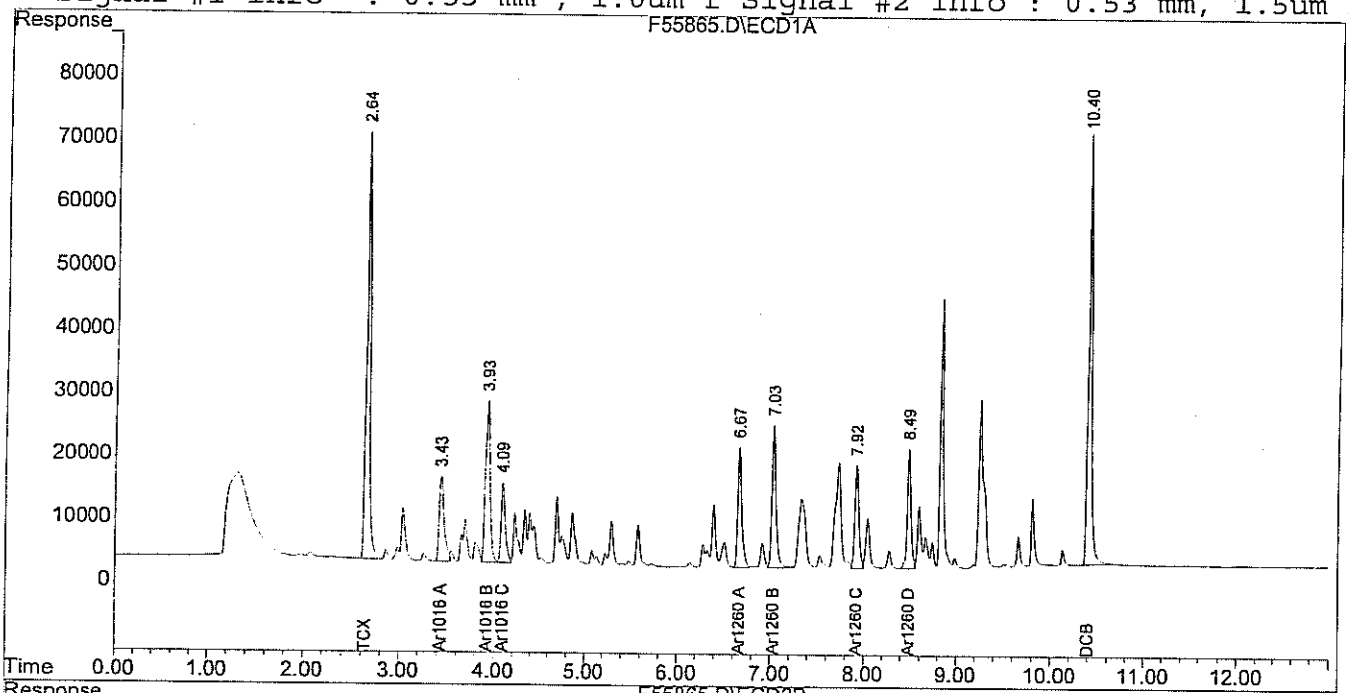
Compound	RT#1	RT#2	Resp#1	Resp#2	ug/ml	ug/ml
System Monitoring Compounds						
1) S TCX	2.64	2.73	1837346	2057065	0.058	0.058
Spiked Amount 0.100	Range 46 - 122		Recovery =		58.00%	58.00%
9) S DCB	10.40	10.04	1671904	2550743	0.054	0.061
Spiked Amount 0.100	Range 26 - 86		Recovery =		54.00%	61.00%
Target Compounds						
2) L1 Ar1016 A	3.44	3.37	469862	514274	0.296	0.300
3) L1 Ar1016 B	3.93	3.87	885341	1150893	0.293	0.299
4) L1 Ar1016 C	4.09	4.03	365795	438601	0.294	0.304
Sum Ar1016 A			1720998	2103768	0.884	0.903
Average Ar1016 A					0.295	0.301
5) L2 Ar1260 A	6.67	6.35	562785	744547	0.288	0.321
6) L2 Ar1260 B	7.03	6.82	670680	1091231	0.285	0.319
7) L2 Ar1260 C	7.92	7.53	510810	774876	0.279	0.318
8) L2 Ar1260 D	8.49	8.59	500733	1688943	0.278	0.299
Sum Ar1260 A			2245008	4299597	1.130	1.256
Average Ar1260 A					0.283	0.314

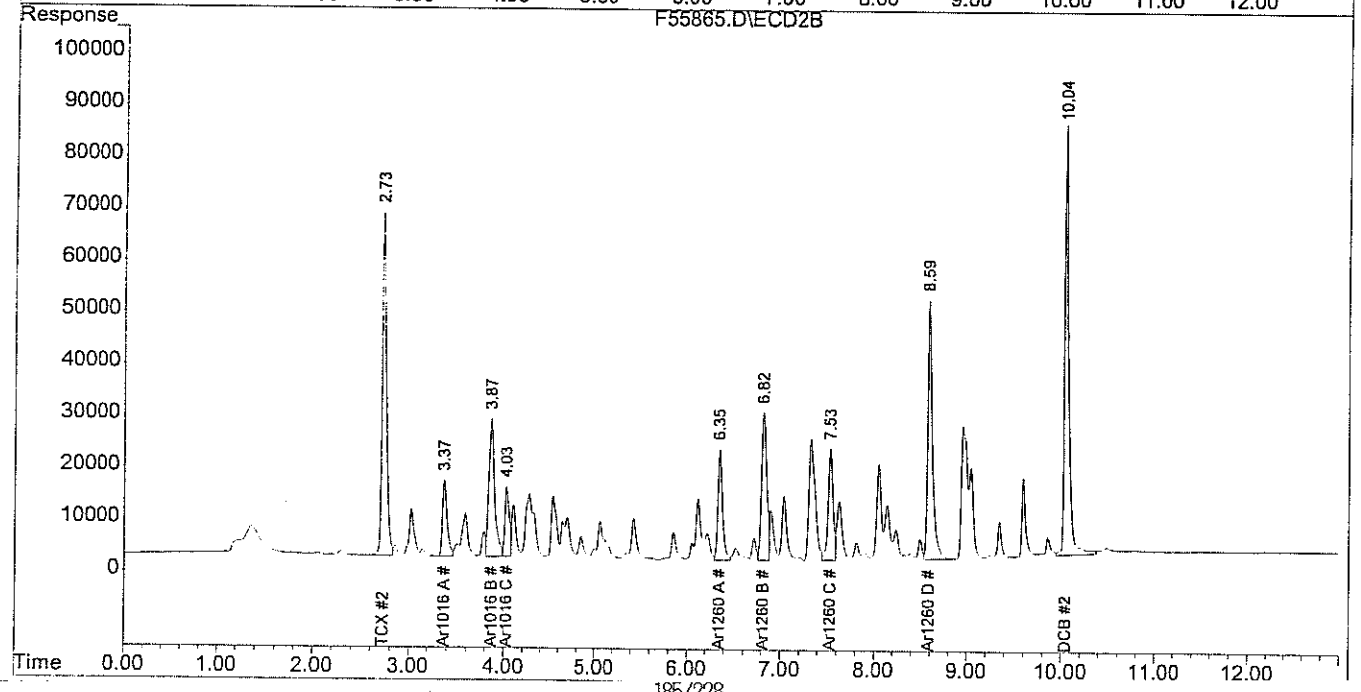
Quantitation Report (QT Reviewed)

Signal #1 : C:\HPCHEM\2\DATA\051606-F\F55865.D\ECD1A.CH Vial: 17
Signal #2 : C:\HPCHEM\2\DATA\051606-F\F55865.D\ECD2B.CH
Acq On : 16 May 2006 8:36 pm Operator:
Sample : P4020 16/60 0.25 Inst : Instrumen
Misc : Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 17 9:54 2006 Quant Results File: PBQ05166.RES

Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 09:54:21 2006
Response via : Multiple Level Calibration
DataAcq Meth : PSTQUICK.M

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film





Quantitation Report (QT Reviewed)

Signal #1 : C:\HPCHEM\2\DATA\051606-F\F55866.D\ECD1A.CH Vial: 18
Signal #2 : C:\HPCHEM\2\DATA\051606-F\F55866.D\ECD2B.CH
Acq On : 16 May 2006 8:51 pm Operator:
Sample : P4021 16/60 0.5 Inst : Instrumen
Misc : Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 17 10:05 2006 Quant Results File: PBQ05166.RES

Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 09:54:21 2006
Response via : Initial Calibration
DataAcq Meth : PSTQUICK.M

5.17.06 5.17.06

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film

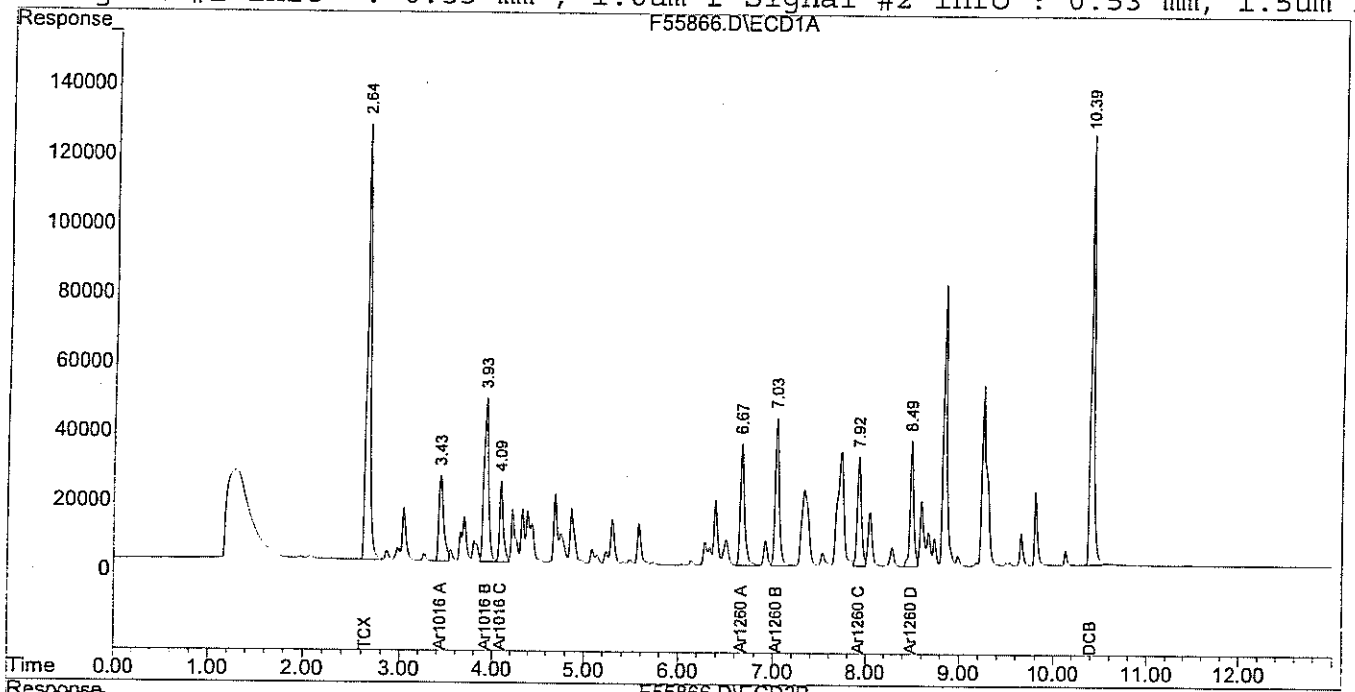
Compound	RT#1	RT#2	Resp#1	Resp#2	ug/ml	ug/ml
System Monitoring Compounds						
1) S TCX	2.64	2.73	3525739	3853418	0.112	0.110
Spiked Amount 0.100	Range 46 - 122		Recovery =		112.00%	110.00%
9) S DCB	10.39	10.04	3091651	4358733	0.104	0.108
Spiked Amount 0.100	Range 26 - 86		Recovery =		104.00%#	108.00%#
Target Compounds						
2) L1 Ar1016 A	3.43	3.37	867015	920329	0.561	0.558
3) L1 Ar1016 B	3.93	3.87	1650244	2072594	0.559	0.557
4) L1 Ar1016 C	4.09	4.03	686360	787671	0.560	0.564
Sum Ar1016 A			3203619	3780595	1.680	1.679
Average Ar1016 A					0.560	0.560
5) L2 Ar1260 A	6.67	6.35	1071560	1271657	0.562	0.575
6) L2 Ar1260 B	7.03	6.82	1268608	1900394	0.552	0.579
7) L2 Ar1260 C	7.92	7.53	993398	1384156	0.549	0.587
8) L2 Ar1260 D	8.49	8.59	972866	3155830	0.547	0.571
Sum Ar1260 A			4306432	7712037	2.210	2.312
Average Ar1260 A					0.553	0.578

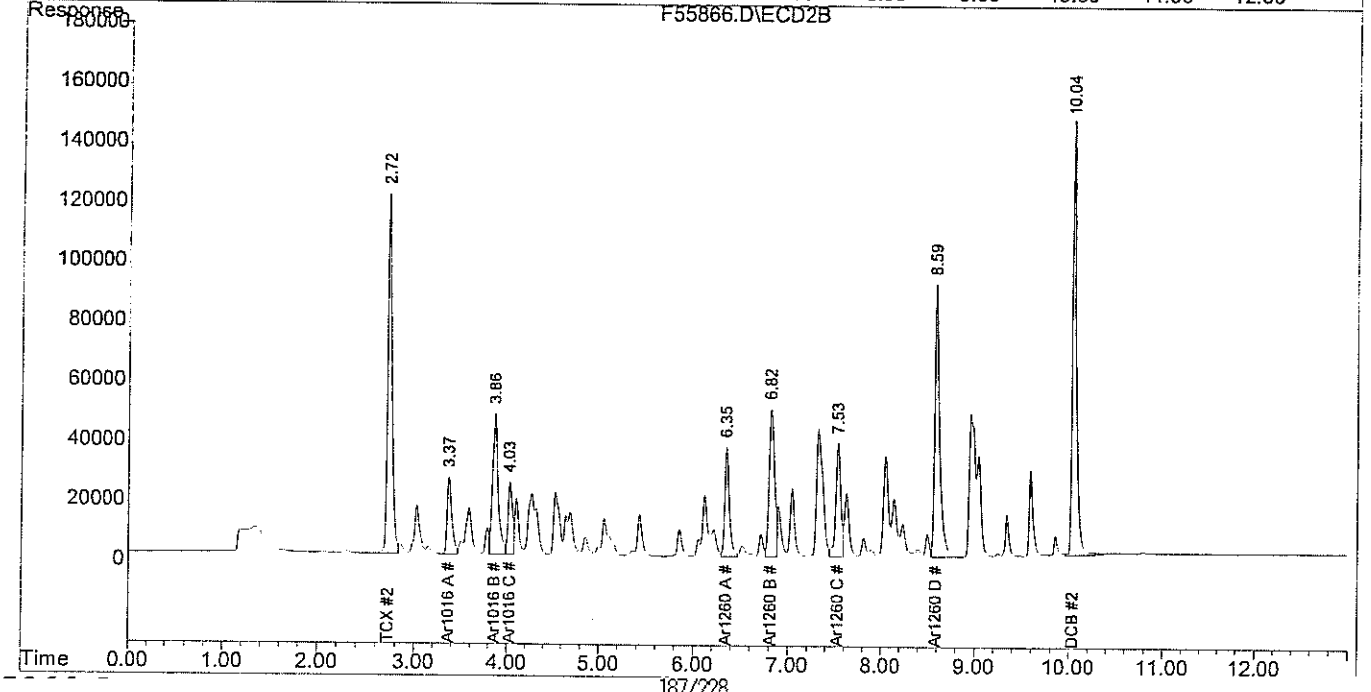
Quantitation Report (QT Reviewed)

Signal #1 : C:\HPCHEM\2\DATA\051606-F\F55866.D\ECD1A.CH Vial: 18
Signal #2 : C:\HPCHEM\2\DATA\051606-F\F55866.D\ECD2B.CH
Acq On : 16 May 2006 8:51 pm Operator:
Sample : P4021 16/60 0.5 Inst : Instrumen
Misc : Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 17 10:05 2006 Quant Results File: PBQ05166.RES

Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 09:54:21 2006
Response via : Multiple Level Calibration
DataAcq Meth : PSTQUICK.M

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film





Quantitation Report (QT Reviewed)

Signal #1 : C:\HPCHEM\2\DATA\051606-F\F55867.D\ECD1A.CH Vial: 19
Signal #2 : C:\HPCHEM\2\DATA\051606-F\F55867.D\ECD2B.CH
Acq On : 16 May 2006 9:20 pm Operator:
Sample : P4022 16/60 1.0 Inst : Instrumen
Misc : Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 17 10:05 2006 Quant Results File: PBQ05166.RES

Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 09:54:21 2006
Response via : Initial Calibration
DataAcq Meth : PSTQUICK.M

5.17.06 5.17.06

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film

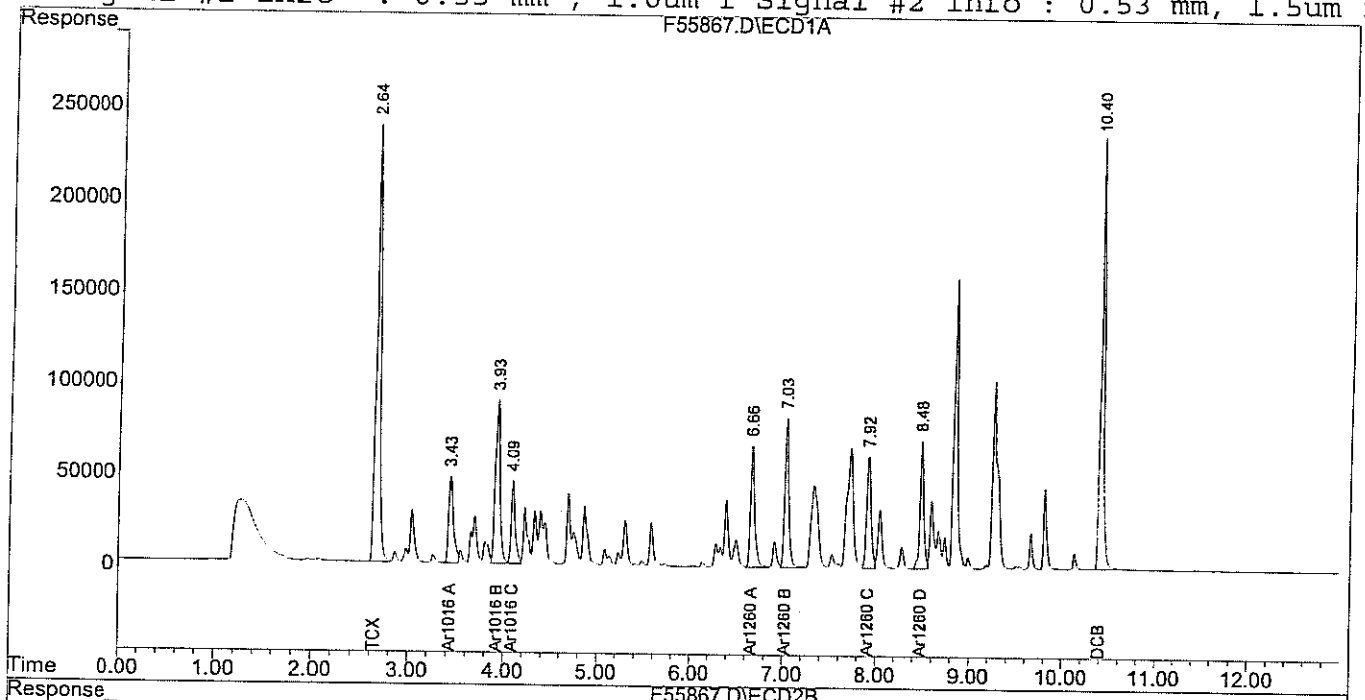
Compound	RT#1	RT#2	Resp#1	Resp#2	ug/ml	ug/ml
System Monitoring Compounds						
1) S TCX	2.64	2.73	6920671	7382493	0.219	0.211
Spiked Amount	0.100	Range 46 - 122		Recovery =	219.00%#	211.00%#
9) S DCB	10.40	10.04	5840708	7809116	0.200	0.200
Spiked Amount	0.100	Range 26 - 86		Recovery =	200.00%#	200.00%#
Target Compounds						
2) L1 Ar1016 A	3.43	3.37	1663419	1708138	1.092	1.059
3) L1 Ar1016 B	3.93	3.87	3193003	3862796	1.095	1.058
4) L1 Ar1016 C	4.09	4.03	1338659	1462474	1.101	1.068
Sum Ar1016 A			6195082	7033407	3.288	3.185
Average Ar1016 A					1.096	1.062
5) L2 Ar1260 A	6.67	6.35	2047135	2313020	1.087	1.076
6) L2 Ar1260 B	7.03	6.82	2456273	3457427	1.083	1.080
7) L2 Ar1260 C	7.92	7.53	1964211	2559137	1.092	1.107
8) L2 Ar1260 D	8.49	8.59	1930282	5977809	1.093	1.094
Sum Ar1260 A			8397902	14307393	4.355	4.357
Average Ar1260 A					1.089	1.089

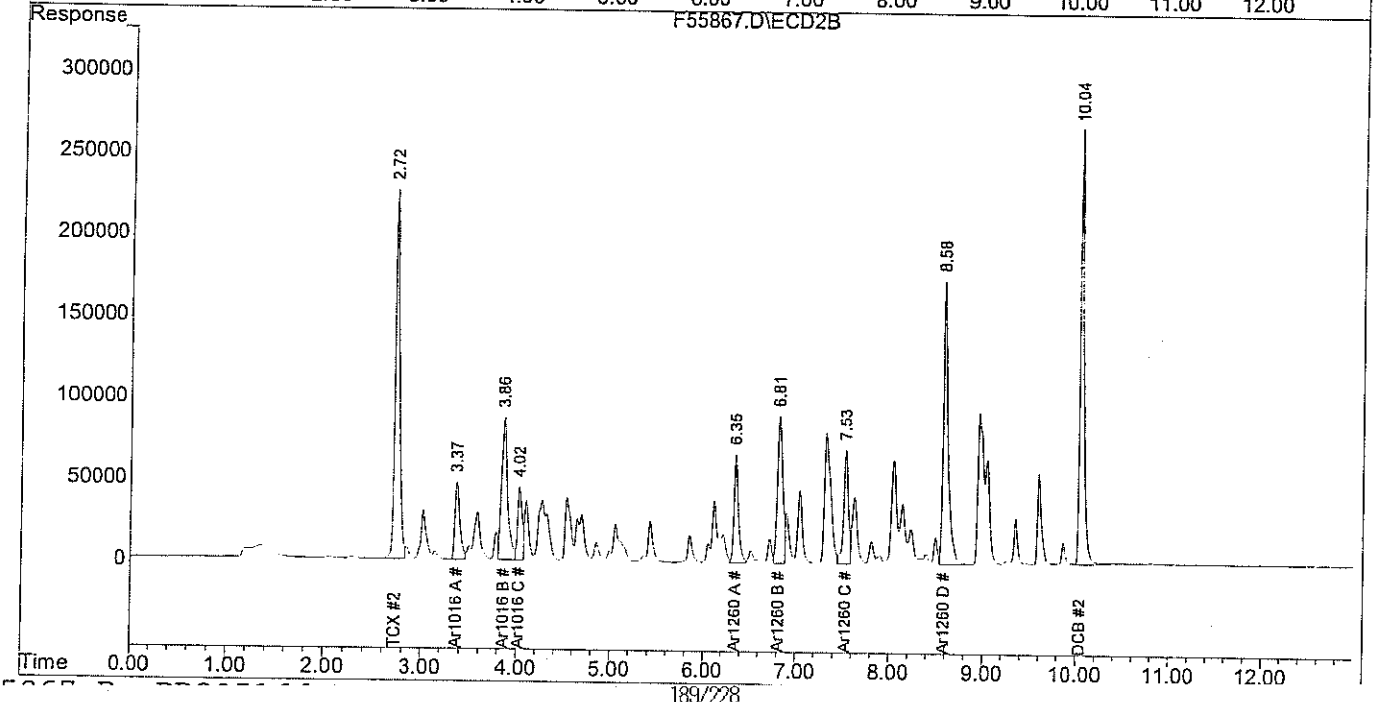
Quantitation Report (QT Reviewed)

Signal #1 : C:\HPCHEM\2\DATA\051606-F\F55867.D\ECD1A.CH Vial: 19
Signal #2 : C:\HPCHEM\2\DATA\051606-F\F55867.D\ECD2B.CH
Acq On : 16 May 2006 9:20 pm Operator:
Sample : P4022 16/60 1.0 Inst : Instrumen
Misc : Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 17 10:05 2006 Quant Results File: PBQ05166.RES

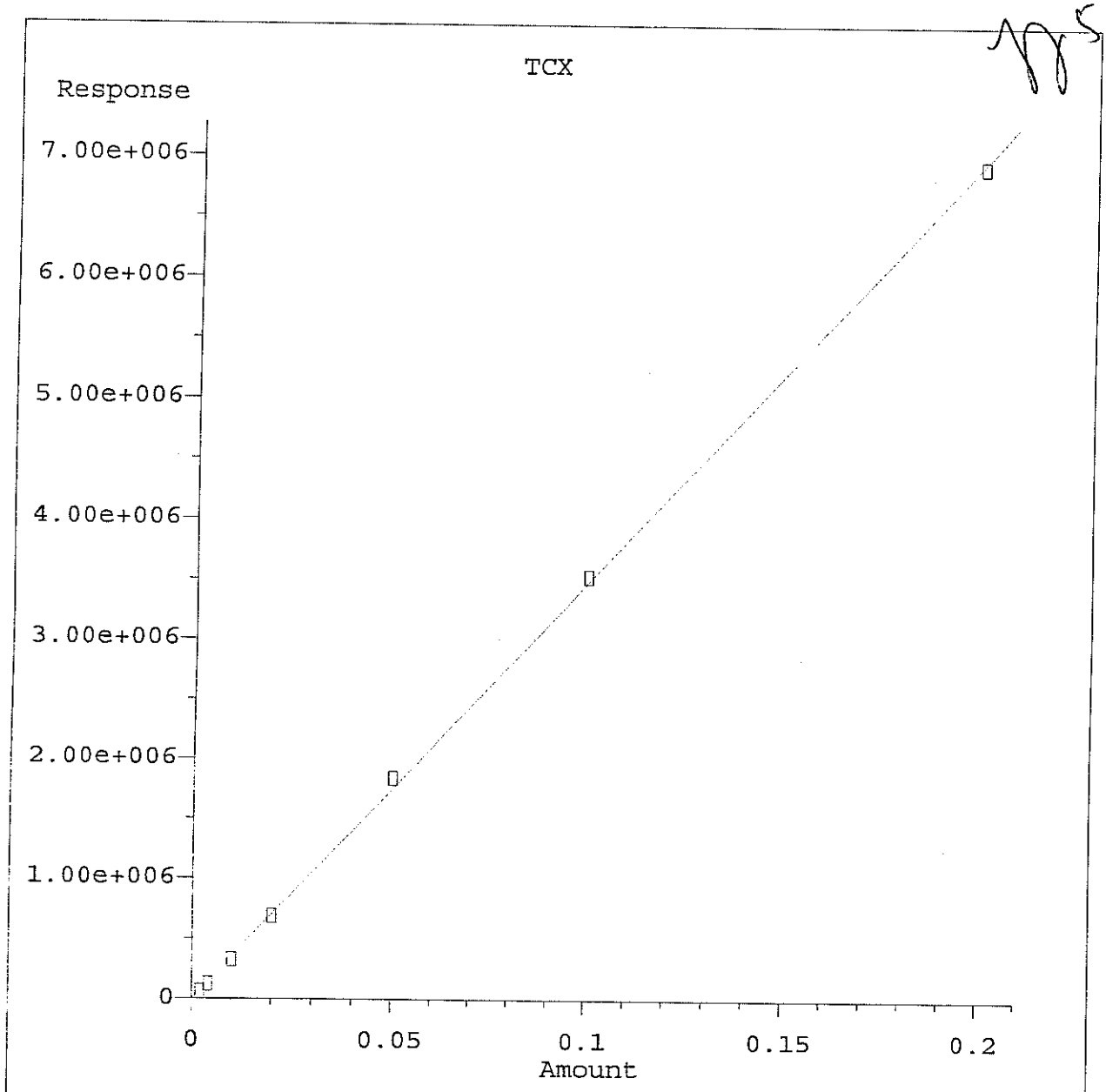
Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 09:54:21 2006
Response via : Multiple Level Calibration
DataAcq Meth : PSTQUICK.M

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film





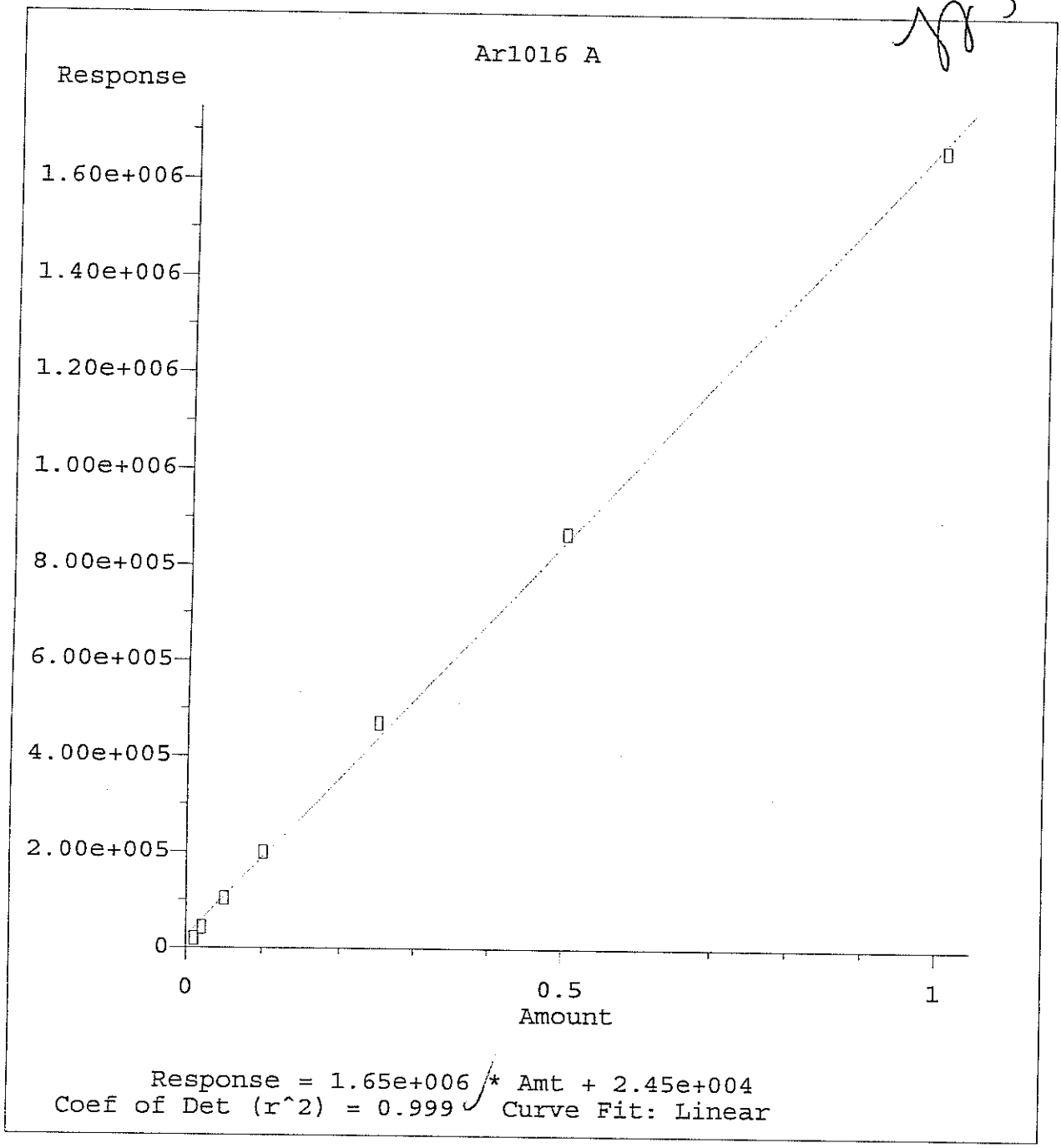
5-17-06



Response = 3.47e+007 * Amt + 1.04e+004
Coef of Det (r^2) = 1.000 Curve Fit: Linear

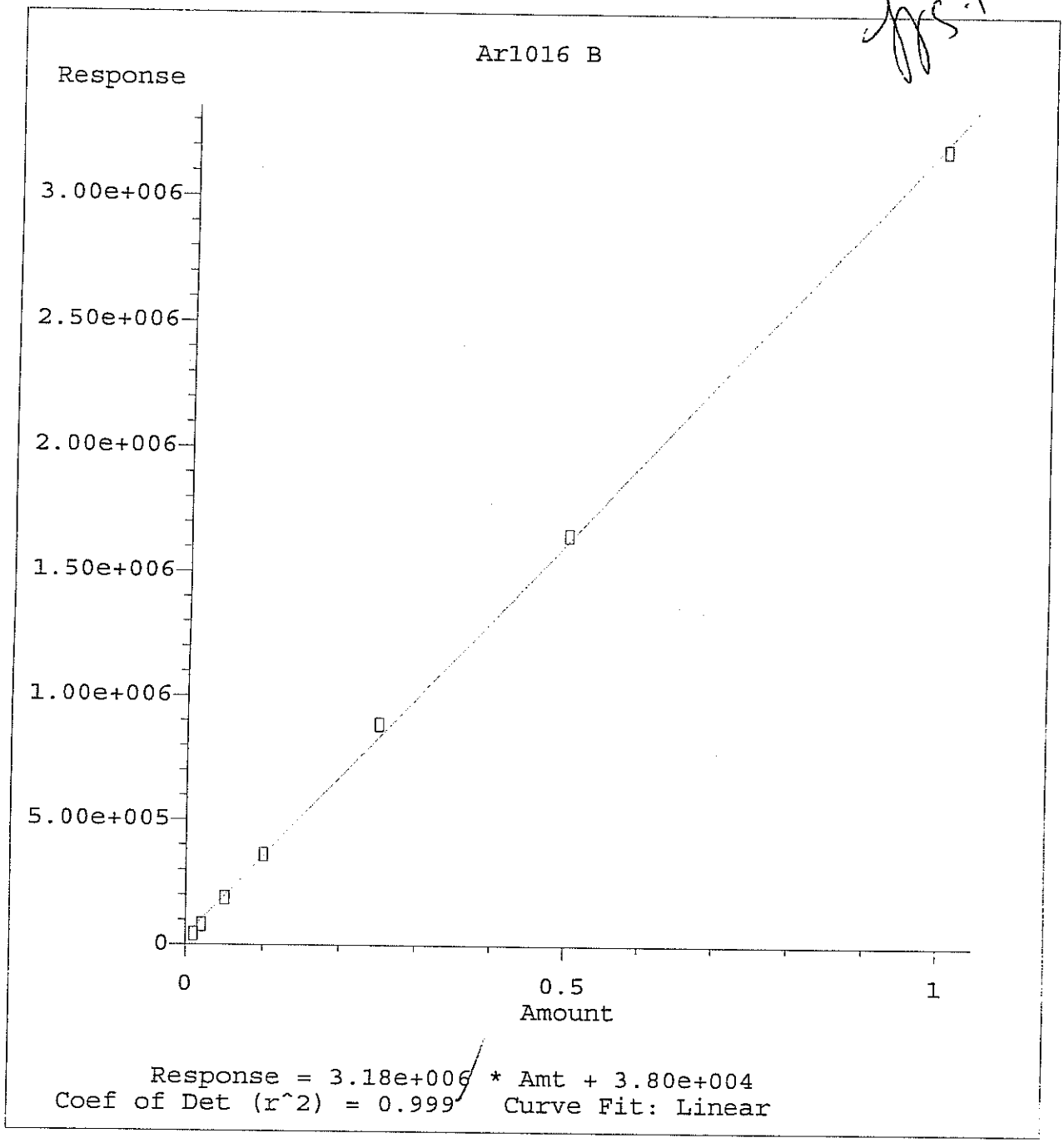
Method Name: C:\HPCHEM\2\METHODS\PBQ05166.M
Calibration Table Last Updated: Wed May 17 10:06:32 2006

5.17.06



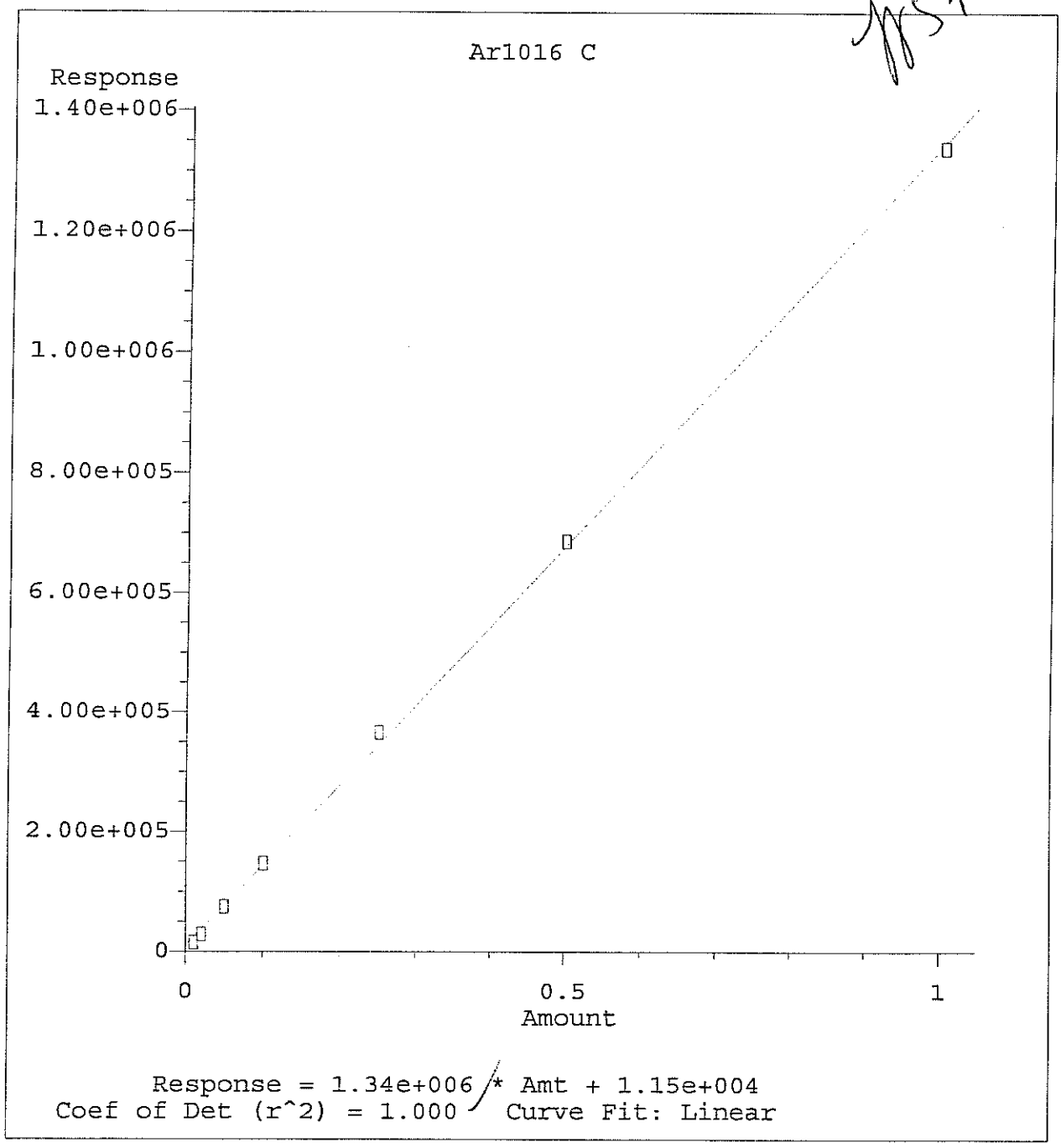
Method Name: C:\HPCHEM\2\METHODS\PBQ05166.M
Calibration Table Last Updated: Wed May 17 10:06:32 2006

5.17.06



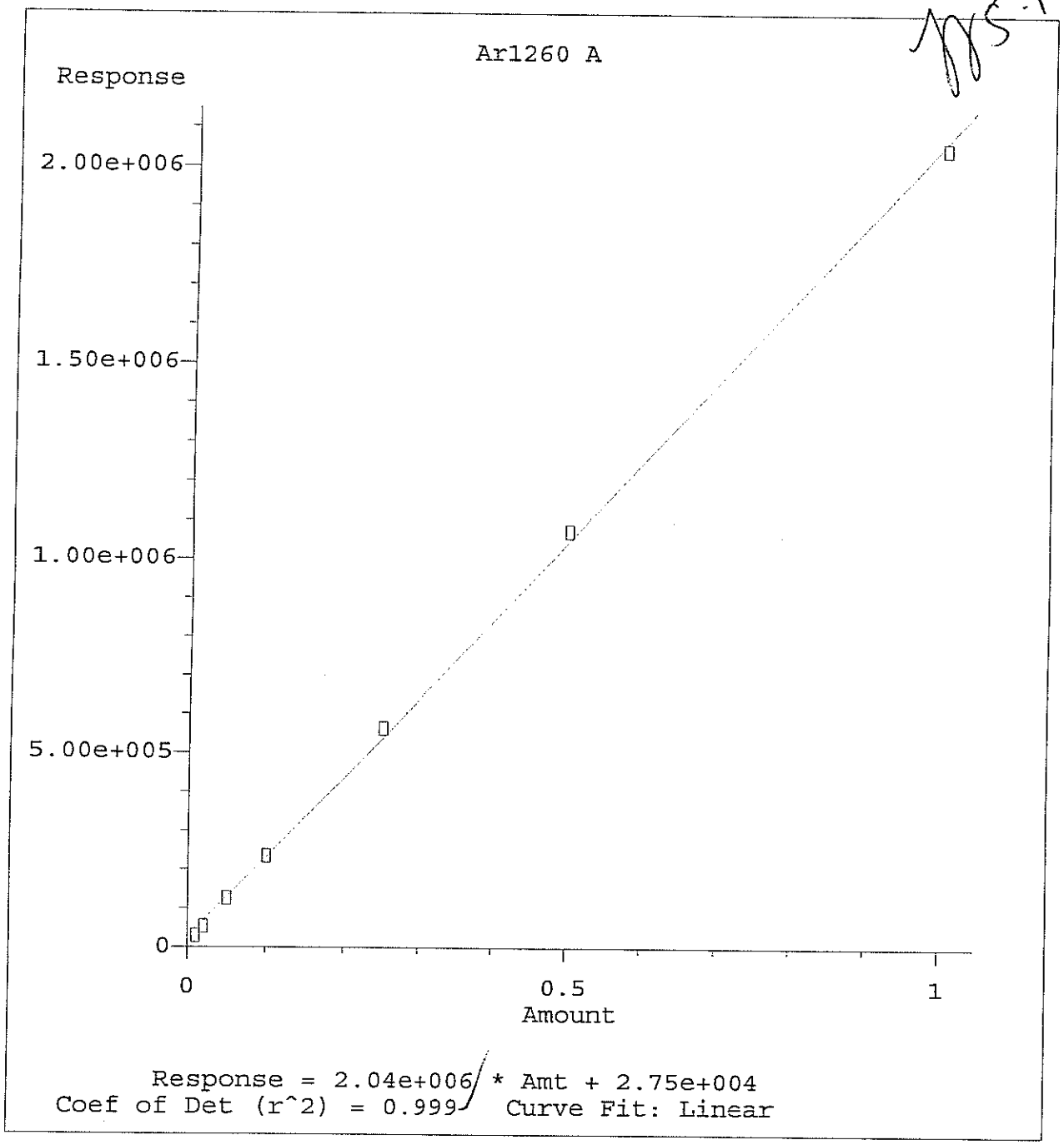
Method Name: C:\HPCHEM\2\METHODS\PBQ05166.M
Calibration Table Last Updated: Wed May 17 10:06:32 2006

5.17.06



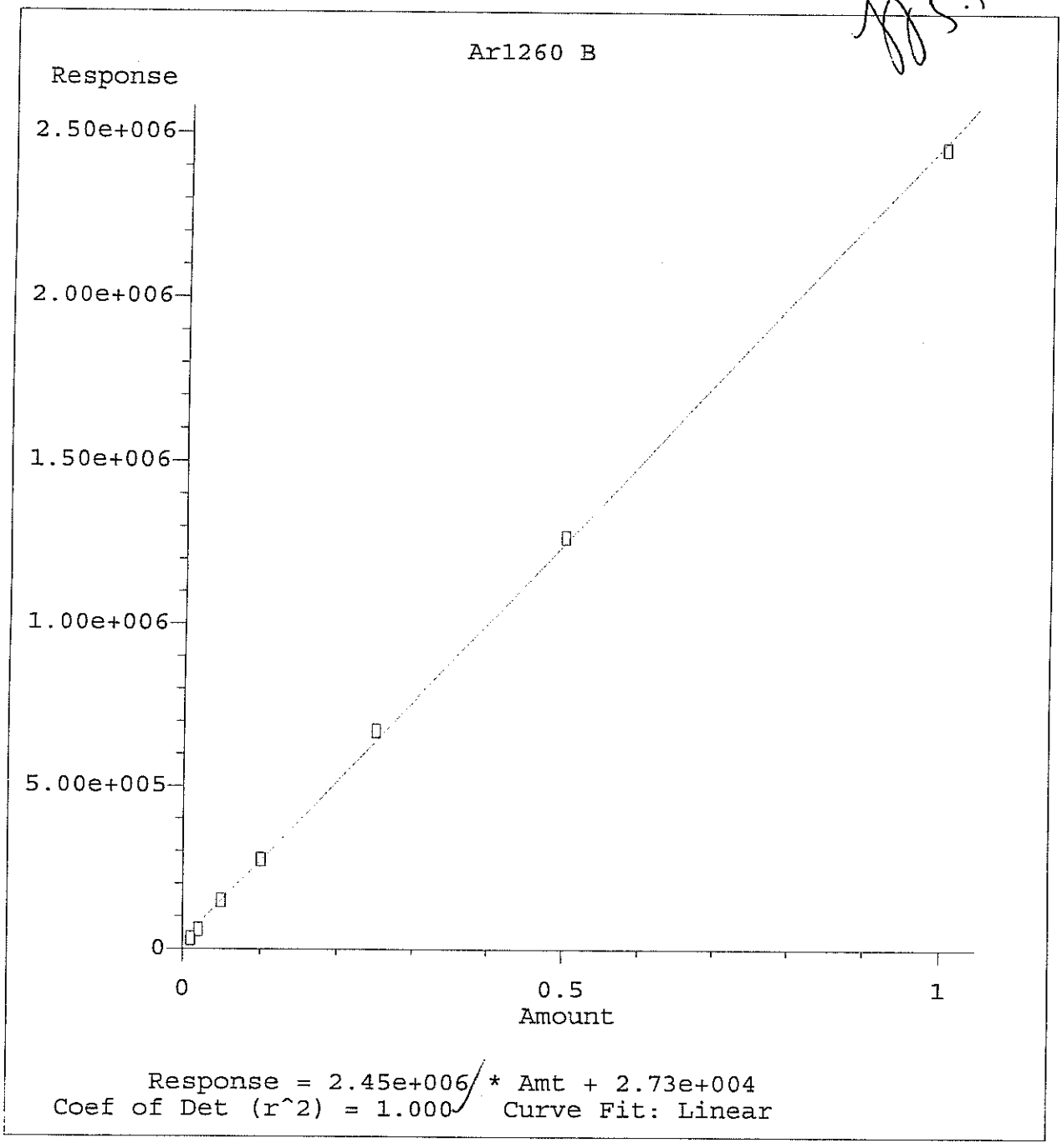
Method Name: C:\HPCHEM\2\METHODS\PBQ05166.M
Calibration Table Last Updated: Wed May 17 10:06:32 2006

JJS-17-06



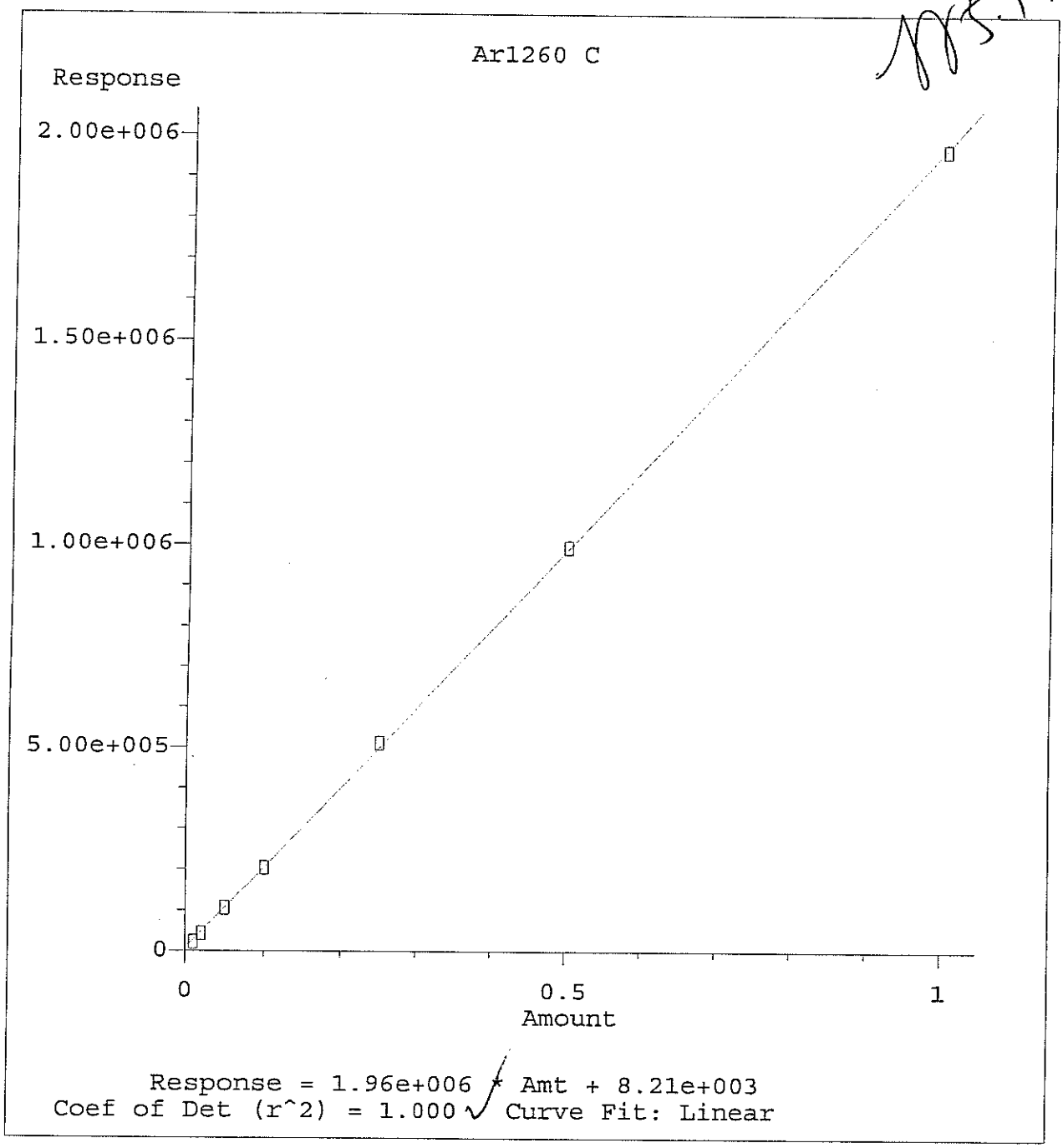
Method Name: C:\HPCHEM\2\METHODS\PBQ05166.M
Calibration Table Last Updated: Wed May 17 10:06:32 2006

5.17.06



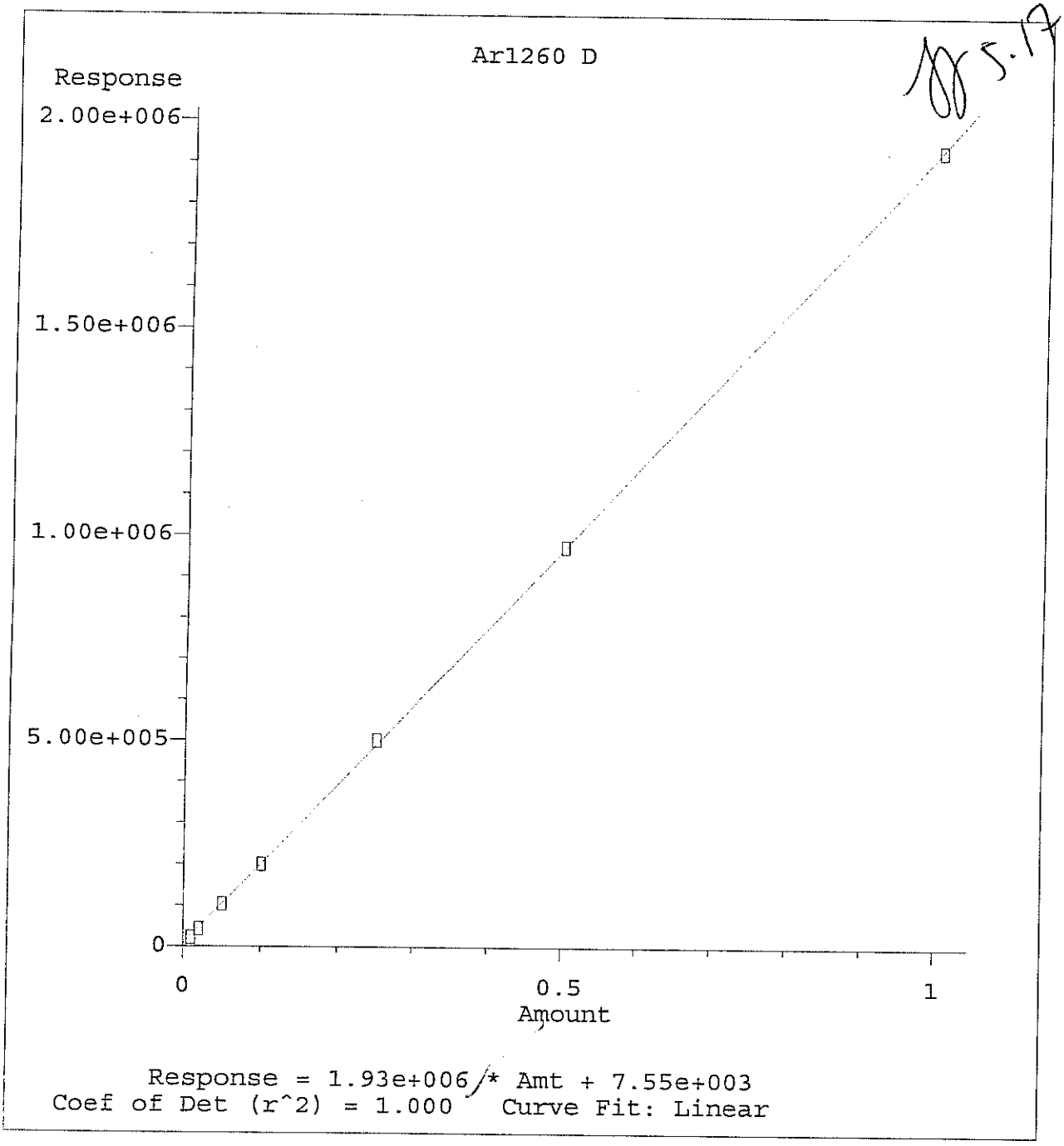
Method Name: C:\HPCHEM\2\METHODS\PBQ05166.M
Calibration Table Last Updated: Wed May 17 10:06:32 2006

5.17.06

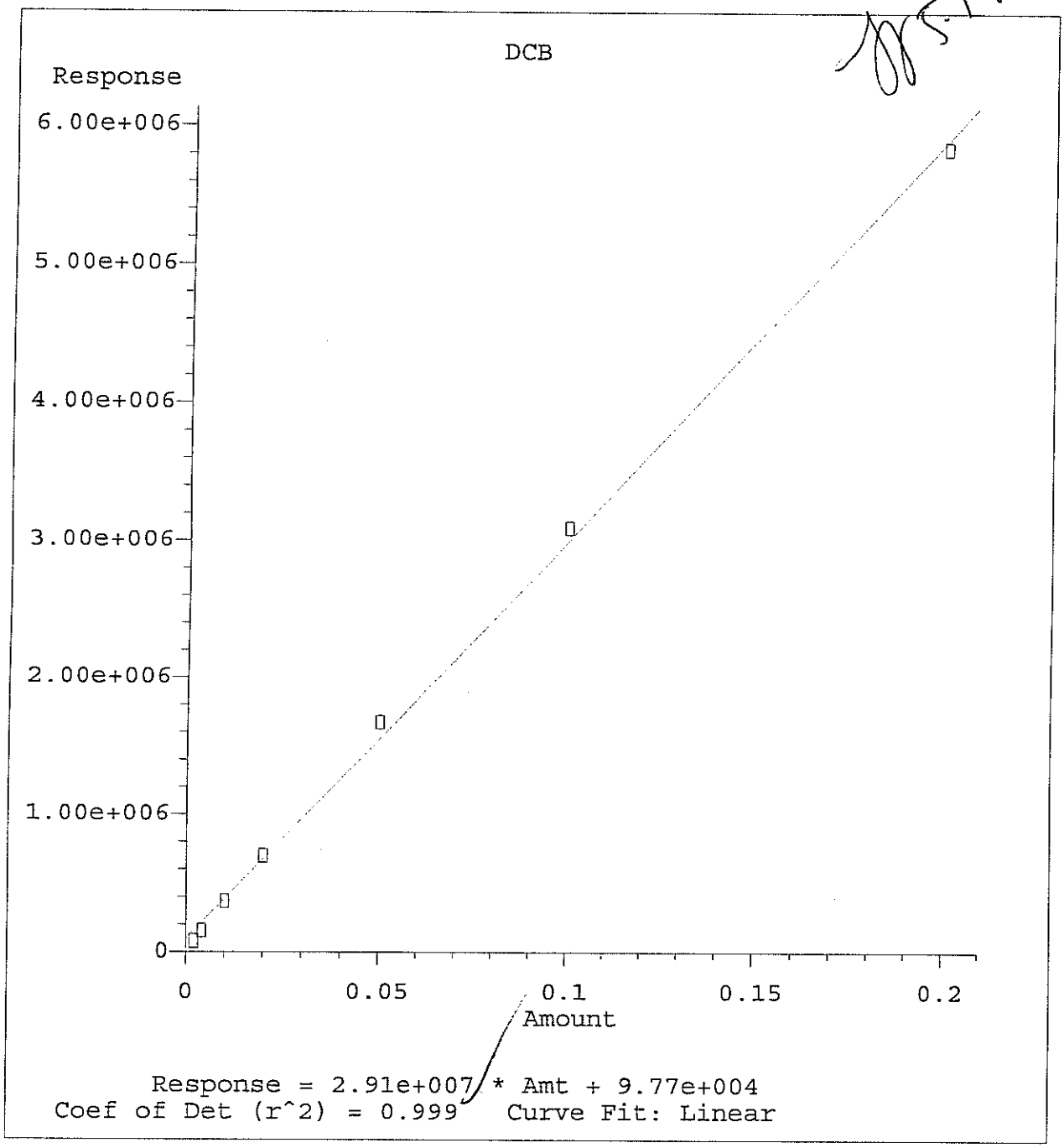


Method Name: C:\HPCHEM\2\METHODS\PBQ05166.M
Calibration Table Last Updated: Wed May 17 10:06:32 2006

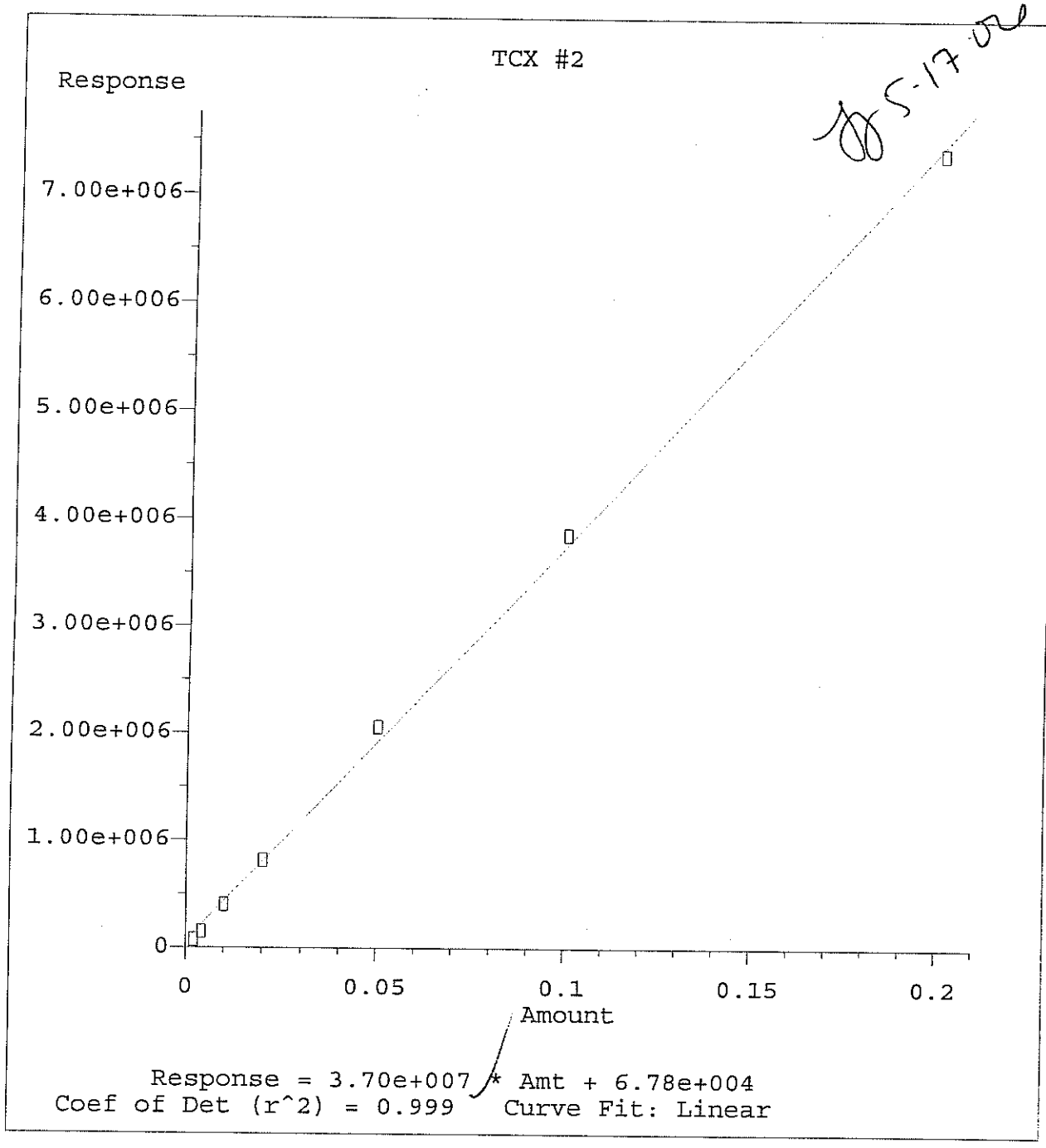
5.17.06



Method Name: C:\HPCHEM\2\METHODS\PBQ05166.M
Calibration Table Last Updated: Wed May 17 10:06:32 2006

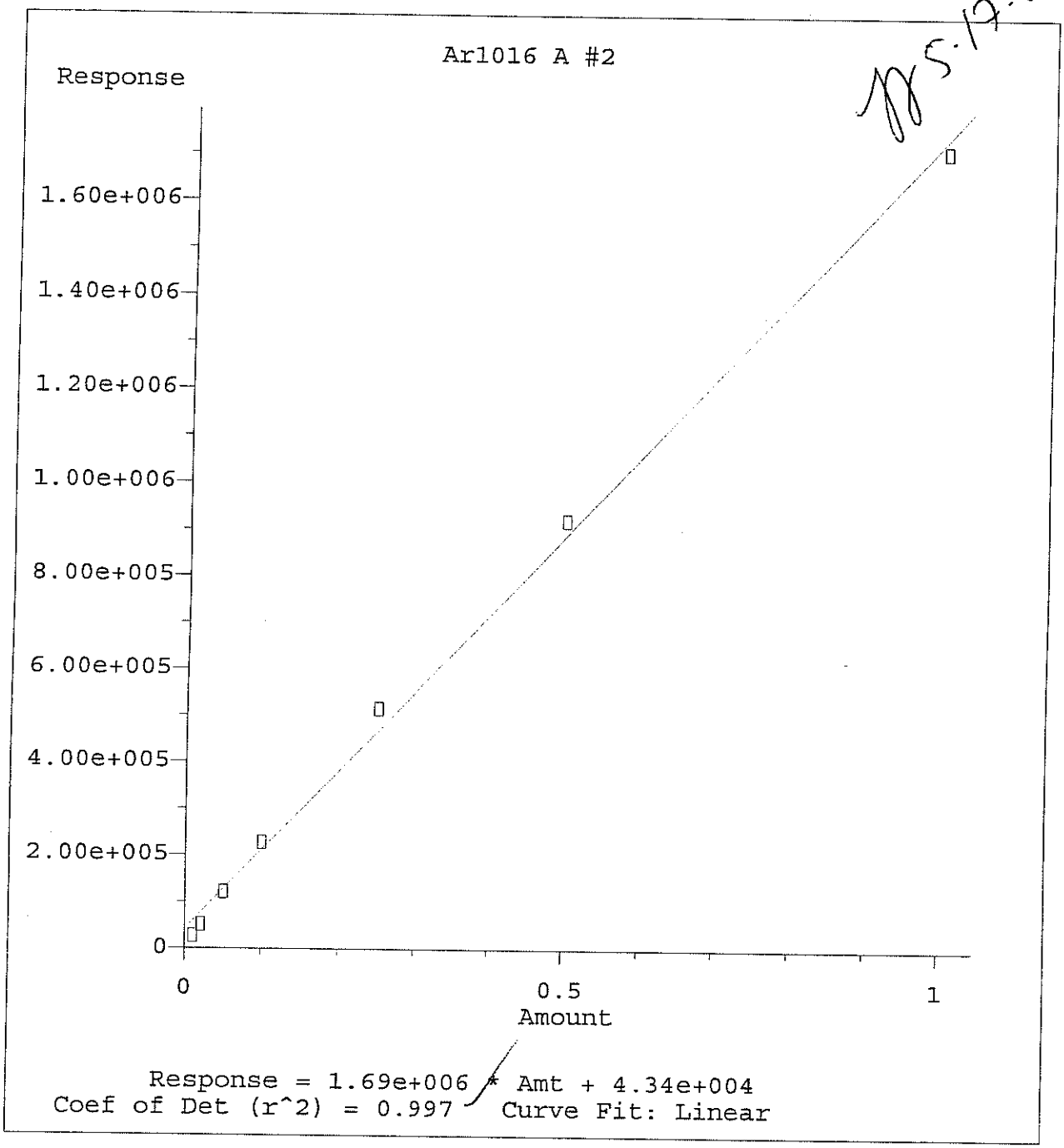


Method Name: C:\HPCHEM\2\METHODS\PBQ05166.M
Calibration Table Last Updated: Wed May 17 10:06:32 2006



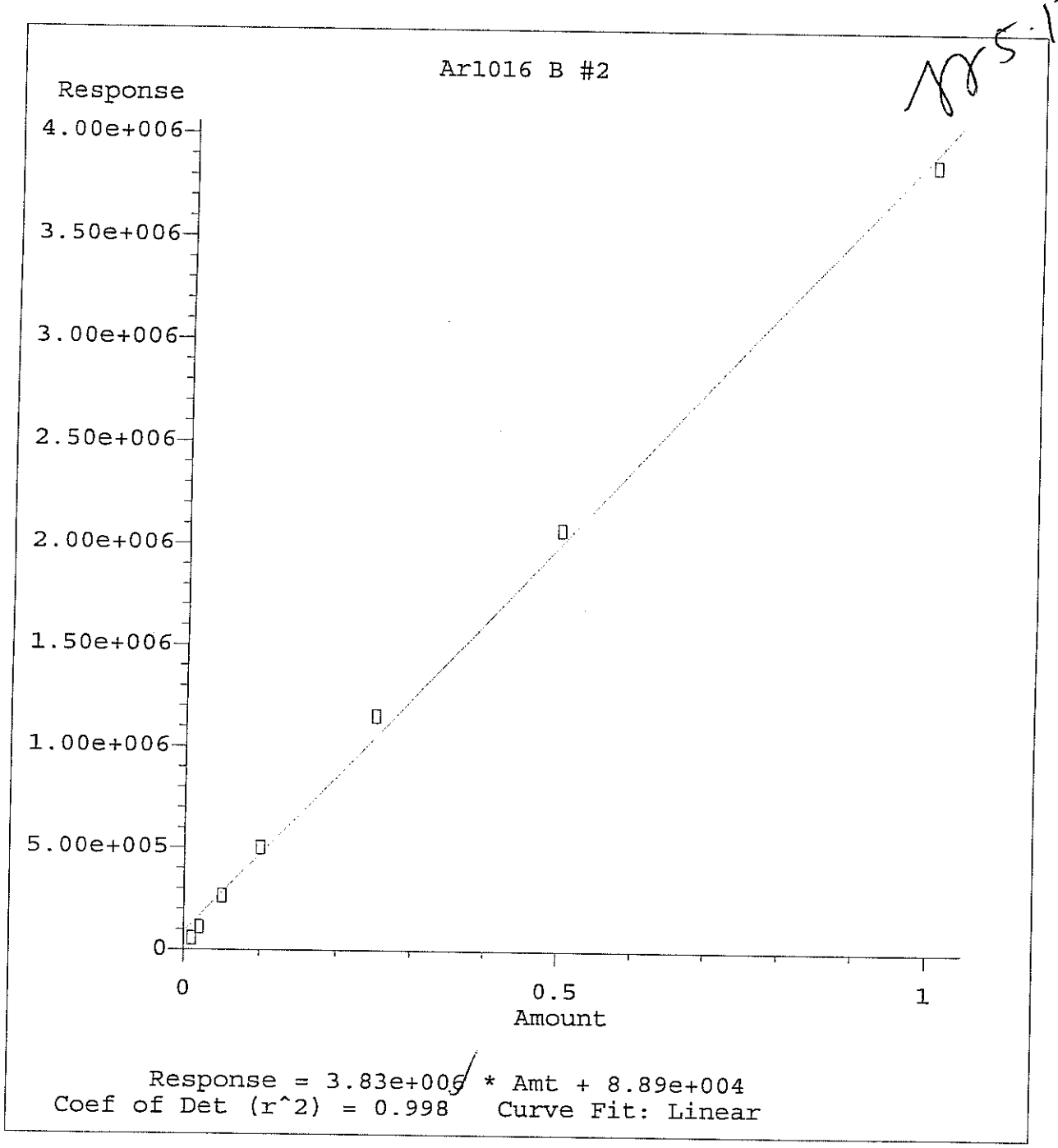
Method Name: C:\HPCHEM\2\METHODS\PBQ05166.M
Calibration Table Last Updated: Wed May 17 10:06:32 2006

5.17.06



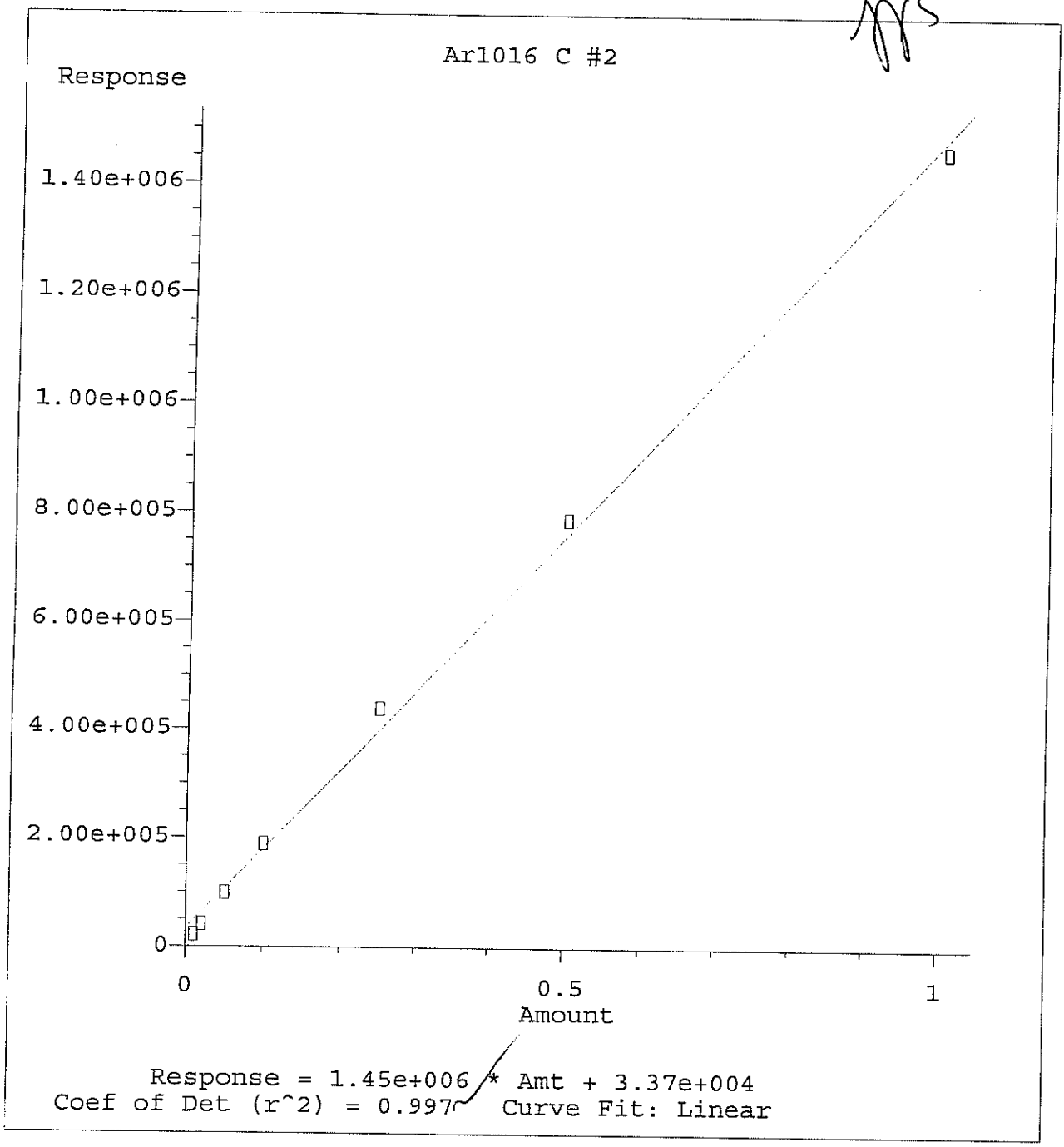
Method Name: C:\HPCHEM\2\METHODS\PBQ05166.M
Calibration Table Last Updated: Wed May 17 10:06:32 2006

5.17.06



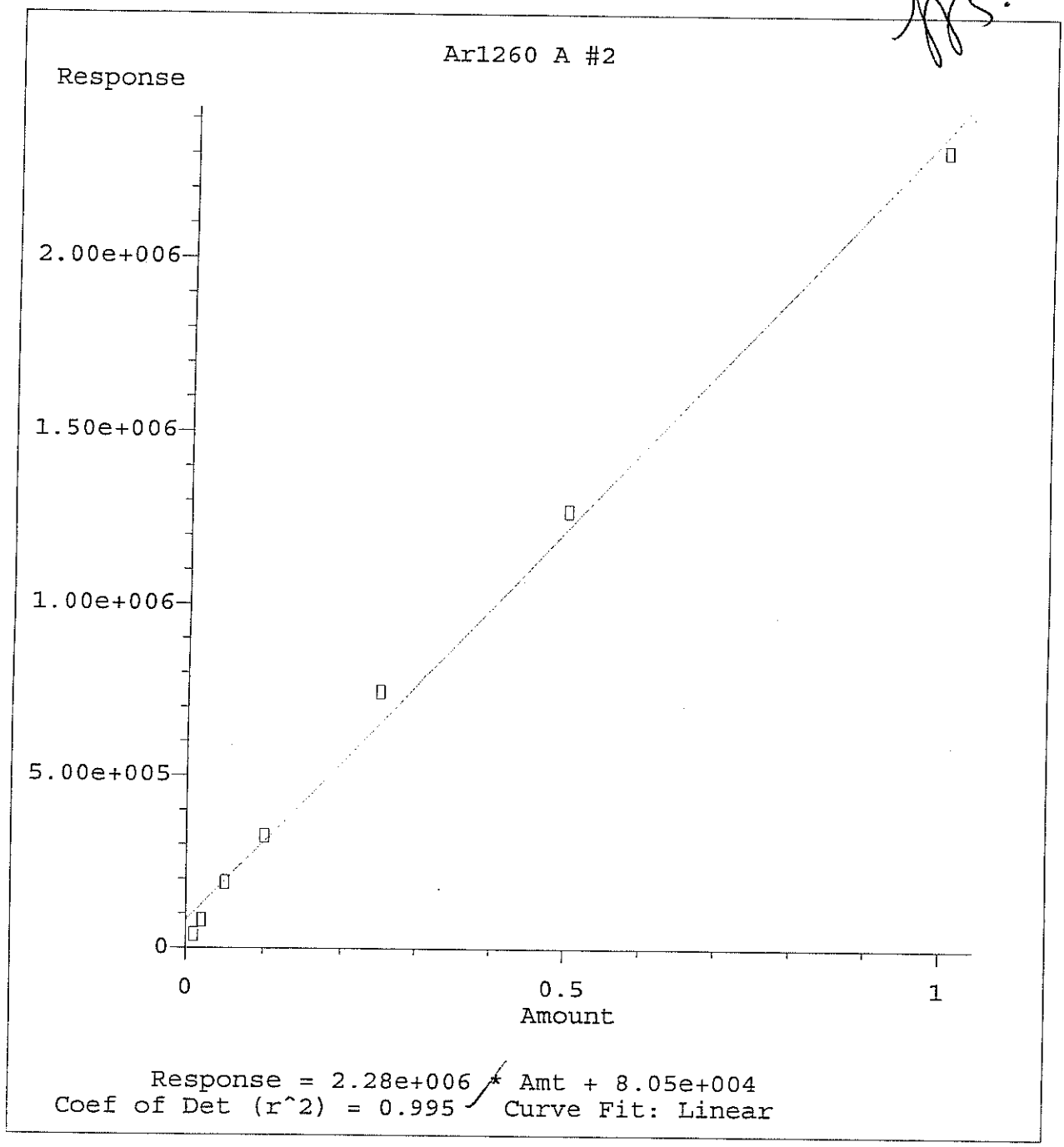
Method Name: C:\HPCHEM\2\METHODS\PBQ05166.M
Calibration Table Last Updated: Wed May 17 10:06:32 2006

5.17.06



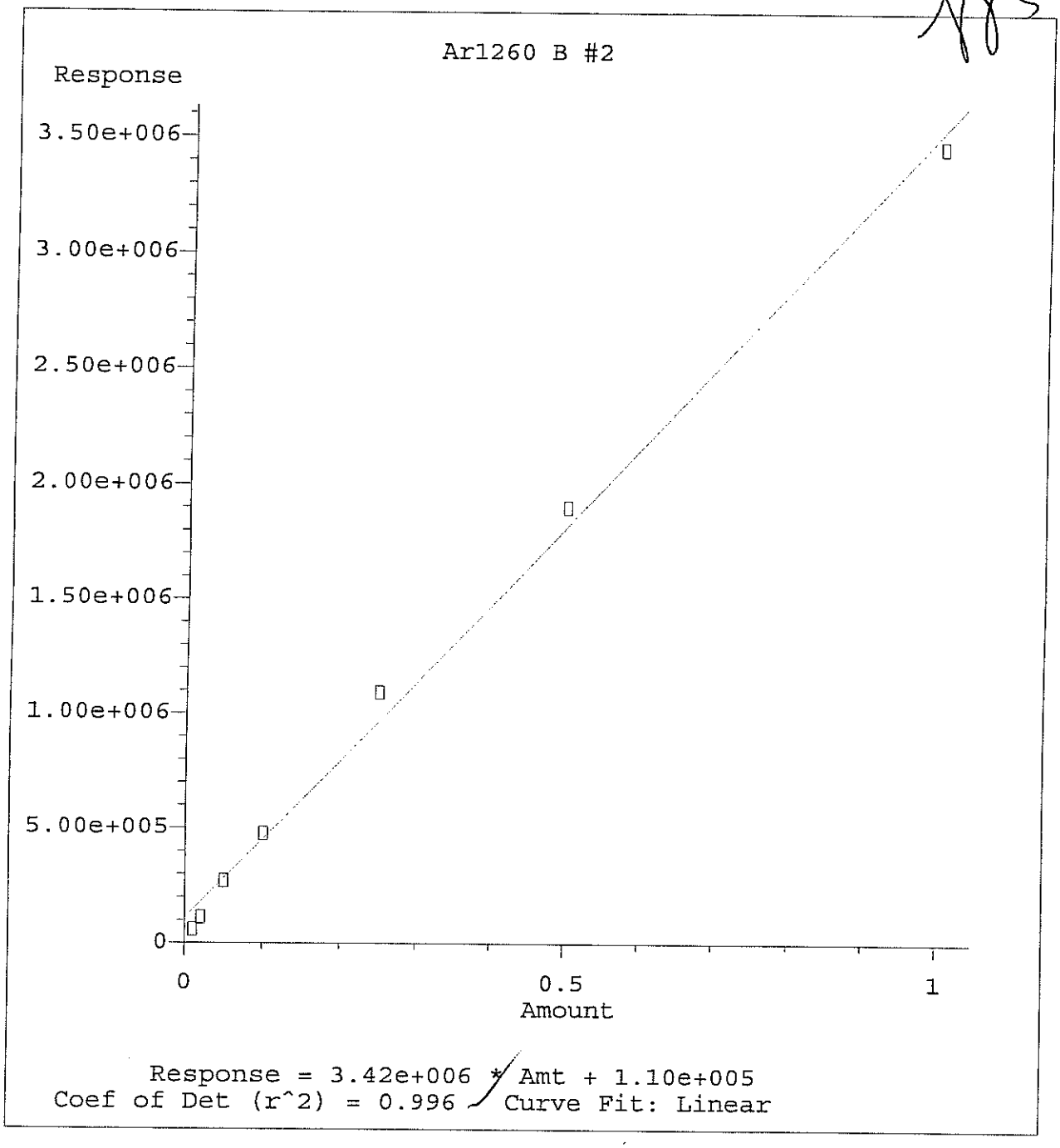
Method Name: C:\HPCHEM\2\METHODS\PBQ05166.M
Calibration Table Last Updated: Wed May 17 10:06:32 2006

5.17.06



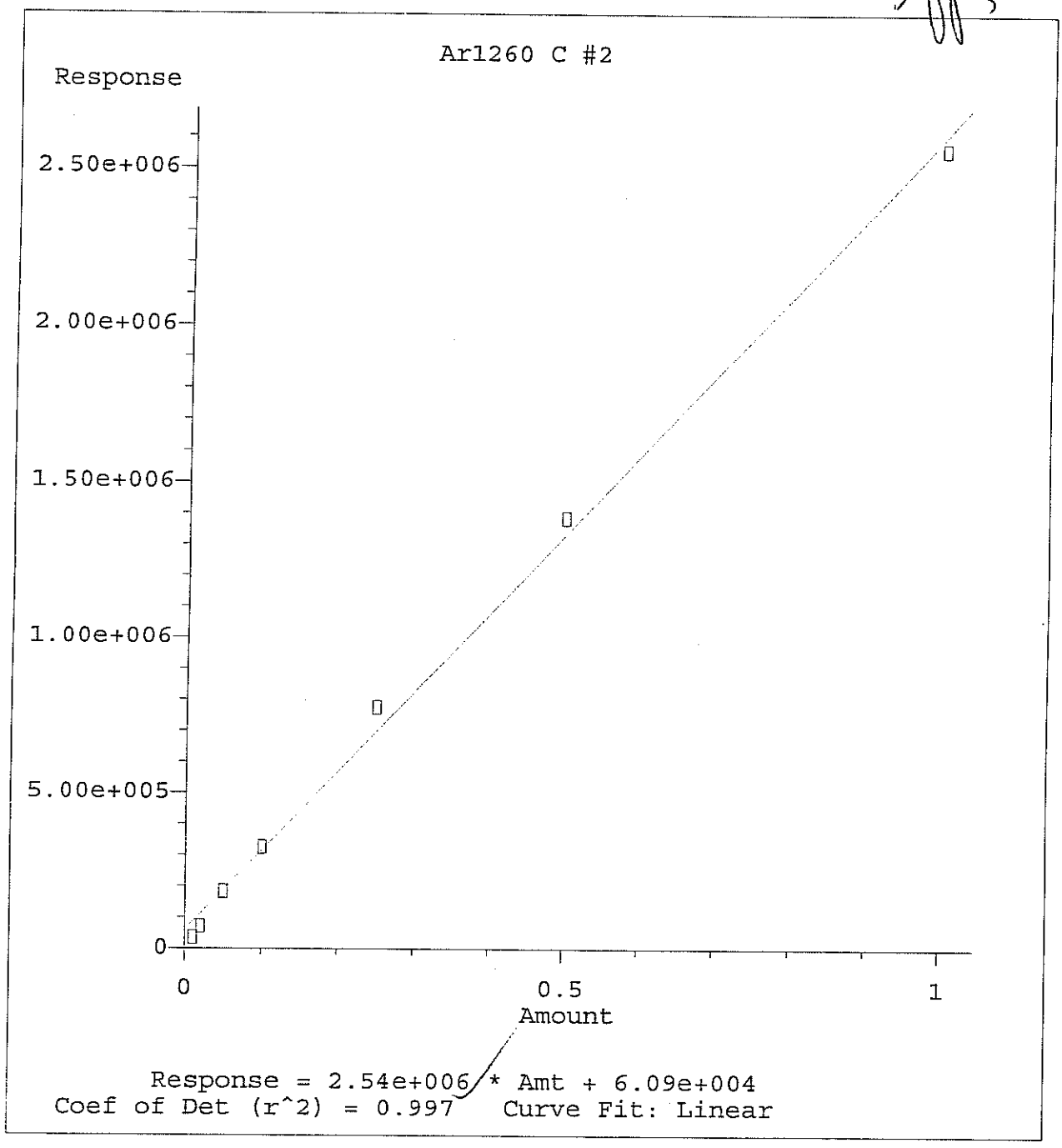
Method Name: C:\HPCHEM\2\METHODS\PBQ05166.M
Calibration Table Last Updated: Wed May 17 10:06:32 2006

5.17.06

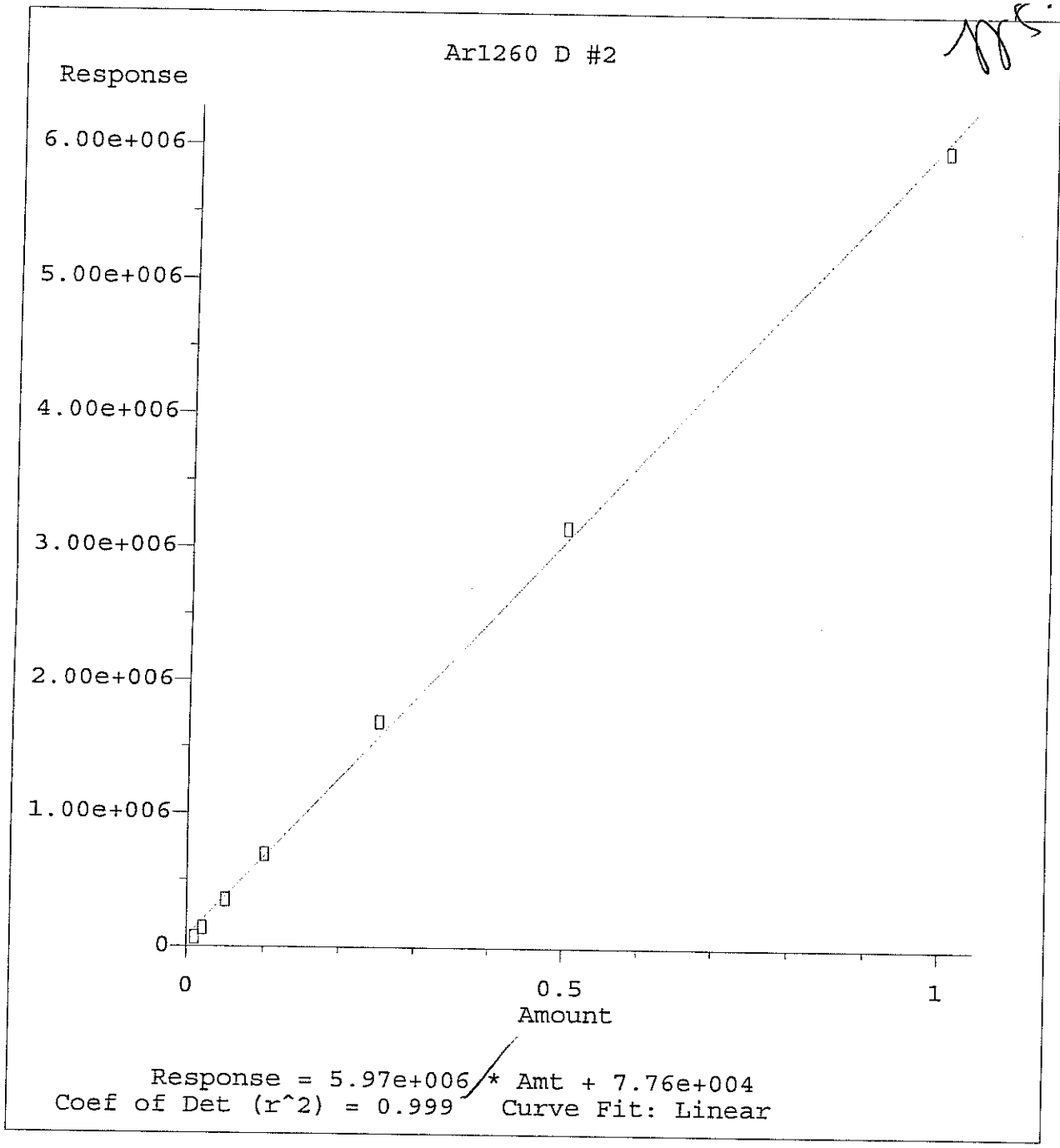


Method Name: C:\HPCHEM\2\METHODS\PBQ05166.M
Calibration Table Last Updated: Wed May 17 10:06:32 2006

5.17.06

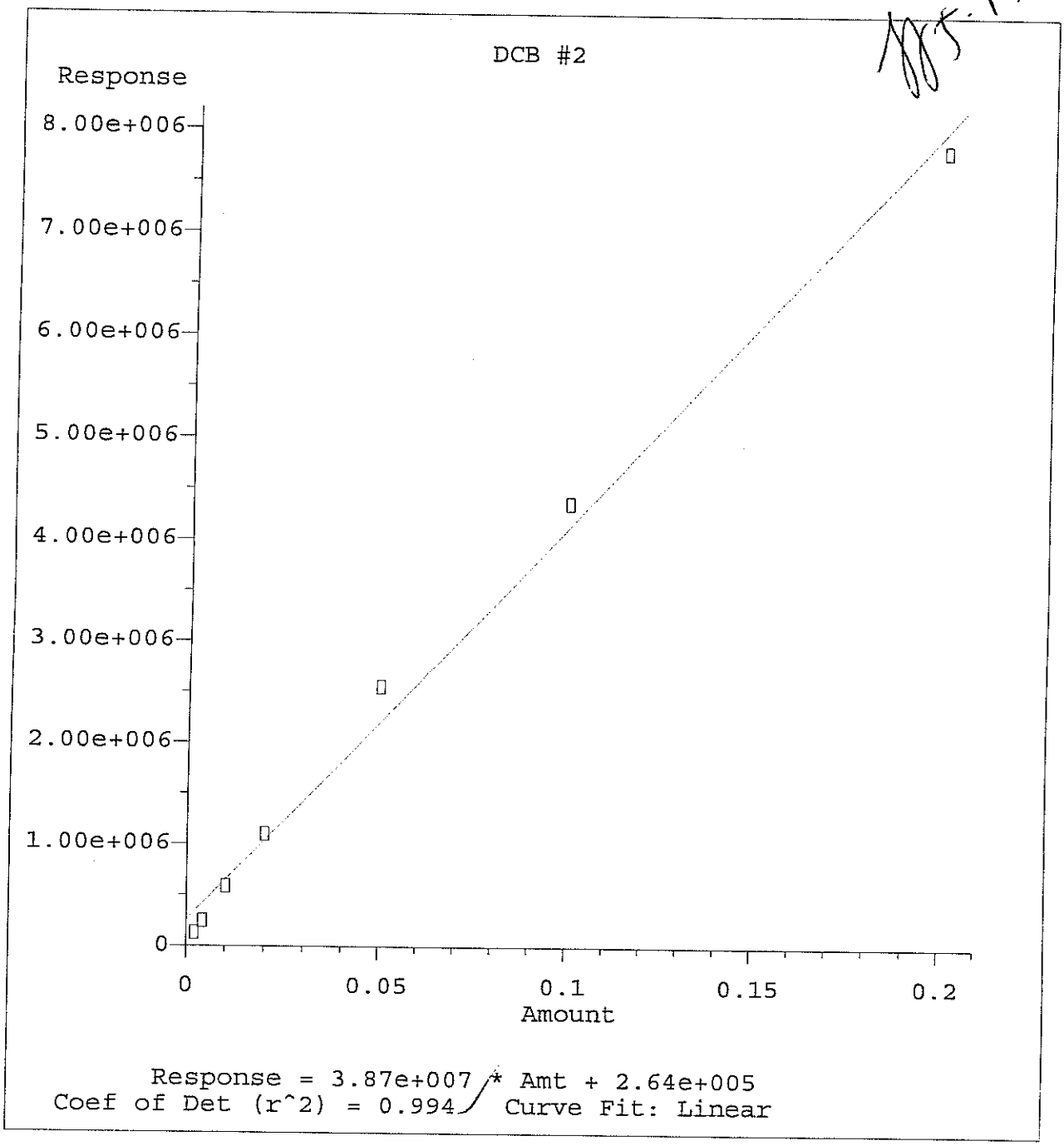


Method Name: C:\HPCHEM\2\METHODS\PBQ05166.M
Calibration Table Last Updated: Wed May 17 10:06:32 2006



Method Name: C:\HPCHEM\2\METHODS\PBQ05166.M
Calibration Table Last Updated: Wed May 17 10:06:32 2006

5.17.06



Method Name: C:\HPCHEM\2\METHODS\PBQ05166.M
Calibration Table Last Updated: Wed May 17 10:06:32 2006

Quantitation Report (Not Reviewed)

Signal #1 : C:\HPCHEM\2\DATA\051606-F\F55868.D\ECD1A.CH Vial: 20
Signal #2 : C:\HPCHEM\2\DATA\051606-F\F55868.D\ECD2B.CH
Acq On : 16 May 2006 9:35 pm Operator:
Sample : P4024 16/60 0.5 CC Inst : Instrumen
Misc : Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 17 10:08 2006 Quant Results File: PBQ05166.RES

Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 10:06:32 2006
Response via : Initial Calibration
DataAcq Meth : PSTQUICK.M

5.17.06 5.17.06

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film

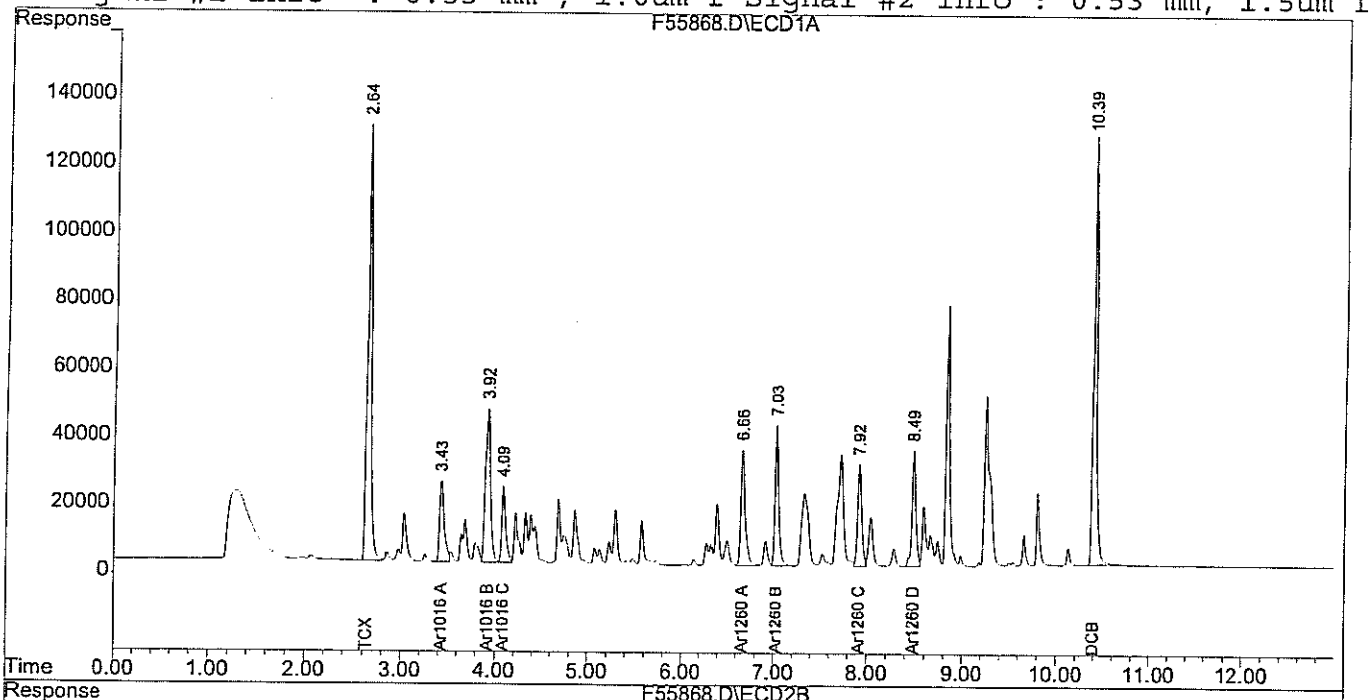
Compound	RT#1	RT#2	Resp#1	Resp#2	ug/ml	ug/ml
System Monitoring Compounds						
1) S TCX	2.64	2.72	3626028	3966484	0.104	0.105
Spiked Amount 0.100	Range 46 - 122			Recovery =	104.00%	105.00%
9) S DCB	10.39	10.04	3115536	4423256	0.104	0.107
Spiked Amount 0.100	Range 26 - 86			Recovery =	104.00%#	107.00%#
Target Compounds						
2) L1 Ar1016 A	3.43	3.37	832389	883397	0.488	0.496
3) L1 Ar1016 B	3.92	3.86	1587624	2003833	0.487	0.500
4) L1 Ar1016 C	4.09	4.02	656707	761021	0.483	0.501
Sum Ar1016 A			3076721	3648252	1.459	1.496
Average Ar1016 A					0.486 97%	0.499 100%
5) L2 Ar1260 A	6.66	6.35	1060994	1236509	0.507	0.507
6) L2 Ar1260 B	7.03	6.82	1233068	1866651	0.493	0.514
7) L2 Ar1260 C	7.92	7.53	954369	1330389	0.482	0.499
8) L2 Ar1260 D	8.49	8.59	934025	2982100	0.481	0.486
Sum Ar1260 A			4182456	7415649	1.963	2.006
Average Ar1260 A					0.491 99%	0.501 100%

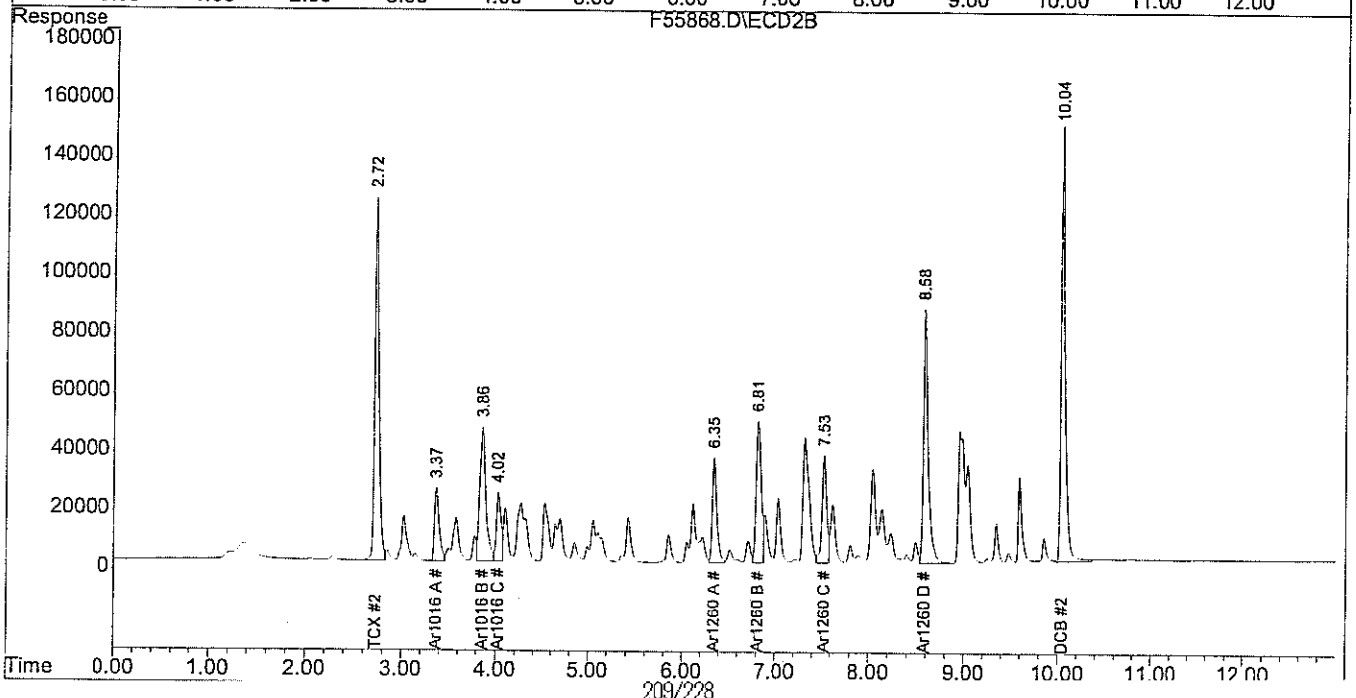
Quantitation Report (Not Reviewed)

Signal #1 : C:\HPCHEM\2\DATA\051606-F\F55868.D\ECD1A.CH Vial: 20
Signal #2 : C:\HPCHEM\2\DATA\051606-F\F55868.D\ECD2B.CH
Acq On : 16 May 2006 9:35 pm Operator:
Sample : P4024 16/60 0.5 CC Inst : Instrumen
Misc : Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 17 10:08 2006 Quant Results File: PBQ05166.RES

Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 10:06:32 2006
Response via : Multiple Level Calibration
DataAcq Meth : PSTQUICK.M

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film





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7A

PCB CONTINUING CALIBRATION CHECK

Site ID: ME

Project No.: 218736 GC Column #1: RTX-CLP Pesticides I Data Package ID: 56617

Instrument ID: F GC Column #2: RTX-CLP Pesticides II ID: 0.53 mm

File #: F56057 Cal. Check Date: 5/24/06 15:30

COMPOUND	True Value (ug/mL)	Measured Value (ug/mL)	% Recovery	*	Limits %R	Total Out
PCB 1016	0.25	0.250	100		85-115	0
PCB1260	0.25	0.245	98		85-115	0
PCB 1016#2	0.25	0.266	106		85-115	0
PCB 1260 #2	0.25	0.272	109		85-115	0

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7A

PCB CONTINUING CALIBRATION CHECK

Site ID: ME

Project No.: 218736 GC Column #1: RTX-CLP Pesticides I Data Package ID: 56617

Instrument ID: F GC Column #2: RTX-CLP Pesticides II ID: 0.53 mm

File #: F56069 Cal. Check Date: 5/24/06 20:08

COMPOUND	True Value (ug/mL)	Measured Value (ug/mL)	% Recovery	*	Limits %R	Total Out
PCB 1016	0.50	0.492	98		85-115	0
PCB1260	0.50	0.494	99		85-115	0
PCB 1016#2	0.50	0.502	100		85-115	0
PCB 1260 #2	0.50	0.522	104		85-115	0

7A

PCB CONTINUING CALIBRATION CHECK

Site ID: ME

Project No.: 218736 GC Column #1: RTX-CLP Pesticides I Data Package ID: 56617

Instrument ID: F GC Column #2: RTX-CLP Pesticides II ID: 0.53 mm

File #: F56078 Cal. Check Date: 5/25/06 14:26

COMPOUND	True Value (ug/mL)	Measured Value (ug/mL)	% Recovery	*	Limits %R	Total Out
PCB 1016	0.25	0.250	100		85-115	0
PCB1260	0.25	0.245	98		85-115	0
PCB 1016#2	0.25	0.260	104		85-115	0
PCB 1260 #2	0.25	0.277	111		85-115	0

7A
PCB CONTINUING CALIBRATION CHECK

Site ID: ME

Project No.: 218736 GC Column #1: RTX-CLP Pesticides I Data Package ID: 56617

Instrument ID: F GC Column #2: RTX-CLP Pesticides II ID: 0.53 mm

File #: F56096 Cal. Check Date: 5/25/06 20:03

COMPOUND	True Value (ug/mL)	Measured Value (ug/mL)	% Recovery	*	Limits %R	Total Out
PCB 1016	0.50	0.508	102		85-115	0
PCB1260	0.50	0.504	101		85-115	0
PCB 1016#2	0.50	0.515	103		85-115	0
PCB 1260 #2	0.50	0.543	109		85-115	0

7A

PCB CONTINUING CALIBRATION CHECK

Site ID: ME

Project No.: 218736 GC Column #1: RTX-CLP Pesticides I Data Package ID: 56617

Instrument ID: F GC Column #2: RTX-CLP Pesticides II ID: 0.53 mm

File #: F56109 Cal. Check Date: 5/25/06 23:27

COMPOUND	True Value (ug/mL)	Measured Value (ug/mL)	% Recovery	*	Limits %R	Total Out
PCB 1016	0.50	0.508	102		85-115	0
PCB1260	0.50	0.510	102		85-115	0
PCB 1016#2	0.50	0.516	103		85-115	0
PCB 1260 #2	0.50	0.543	109		85-115	0

7A
PCB CONTINUING CALIBRATION CHECK

Site ID: ME

Project No.: 218736 GC Column #1: RTX-CLP Pesticides I Data Package ID: 56617

Instrument ID: F GC Column #2: RTX-CLP Pesticides II ID: 0.53 mm

File #: F56114 Cal. Check Date: 5/26/06 0:55

COMPOUND	True Value (ug/mL)	Measured Value (ug/mL)	% Recovery	*	Limits %R	Total Out
PCB 1016	0.50	0.497	99		85-115	0
PCB1260	0.50	0.492	98		85-115	0
PCB 1016#2	0.50	0.510	102		85-115	0
PCB 1260 #2	0.50	0.517	103		85-115	0

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7A

PCB CONTINUING CALIBRATION CHECK

Site ID: ME

Project No.: 218736 GC Column #1: RTX-CLP Pesticides I Data Package ID: 56617

Instrument ID: F GC Column #2: RTX-CLP Pesticides II ID: 0.53 mm

File #: F56130 Cal. Check Date: 5/26/06 11:41

COMPOUND	True Value (ug/mL)	Measured Value (ug/mL)	% Recovery	*	Limits %R	Total Out
PCB 1016	0.50	0.538	108		85-115	0
PCB1260	0.50	0.532	106		85-115	0
PCB 1016#2	0.50	0.533	107		85-115	0
PCB 1260 #2	0.50	0.548	110		85-115	0

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7A

PCB CONTINUING CALIBRATION CHECK

Site ID: ME

Project No.: 218736 GC Column #1: RTX-CLP Pesticides I Data Package ID: 56617

Instrument ID: F GC Column #2: RTX-CLP Pesticides II ID: 0.53 mm

File #: F56132 Cal. Check Date: 5/26/06 13:38

COMPOUND	True Value (ug/mL)	Measured Value (ug/mL)	% Recovery	*	Limits %R	Total Out
PCB 1016	0.25	0.277	111		85-115	0
PCB1260	0.25	0.276	110		85-115	0
PCB 1016#2	0.25	0.278	111		85-115	0
PCB 1260 #2	0.25	0.290	116	*	85-115	1

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7A
PCB CONTINUING CALIBRATION CHECK

Site ID: ME

Project No.: 218736 GC Column #1: RTX-CLP Pesticides I Data Package ID: 56617

Instrument ID: F GC Column #2: RTX-CLP Pesticides II ID: 0.53 mm

File #: F56139 Cal. Check Date: 5/26/06 15:49

COMPOUND	True Value (ug/mL)	Measured Value (ug/mL)	% Recovery	*	Limits %R	Total Out
PCB 1016	0.50	0.527	105		85-115	0
PCB1260	0.50	0.532	106		85-115	0
PCB 1016#2	0.50	0.512	102		85-115	0
PCB 1260 #2	0.50	0.559	112		85-115	0

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PCB SYSTEM MONITORING COMPOUND RETENTION TIME SUMMARY

Quote#: SDG: 56617

Project No: 218736 Site ID: Location: ME

GC Columns: RTX-CLPesticidest ID: 0.32(mm)

Instrument ID: F RTX-CLPesticidesII

Date Analyzed: 5/16/06

Lab File ID (Standard) : F55866 Time Analyzed: 20:51

					COLUMN #1		COLUMN #1	
					SMC1	SMC 2	SMC1	SMC 2
	PESTICIDE STANDARD				2.64	10.39	2.73	10.04
	UPPER LIMIT				2.69	10.44	2.78	10.09
	LOWER LIMIT				2.59	10.34	2.68	9.99
	SAMPLE ID	LAB FILE ID	DATE	TIME	COLUMN #1 SMC1 #	COLUMN #1 SMC 2 #	COLUMN #2 SMC1 #	COLUMN #2 SMC 2 #
01	P4016 16/60 0.01	F55861.D	05/16/06	19:38	2.64	10.39	2.72	10.04
02	P4017 16/60 0.02	F55862.D	05/16/06	19:53	2.64	10.40	2.73	10.04
03	P4018 16/60 0.05	F55863.D	05/16/06	20:07	2.64	10.40	2.73	10.05
04	P4019 16/60 0.1	F55864.D	05/16/06	20:22	2.64	10.40	2.73	10.05
05	P4020 16/60 0.25	F55865.D	05/16/06	20:36	2.64	10.40	2.73	10.04
06	P4021 16/60 0.5	F55866.D	05/16/06	20:51	2.64	10.39	2.73	10.04
07	P4022 16/60 1.0	F55867.D	05/16/06	21:20	2.64	10.40	2.73	10.04
08	P4024 16/60 0.5 CC	F55868.D	05/16/06	21:35	2.64	10.39	2.72	10.04
09	P3883 AR 1221	F55869.D	05/16/06	21:50	NA	NA	NA	NA
10	P3884 AR 1232	F55870.D	05/16/06	22:04	NA	NA	NA	NA
11	P3885 AR 1242	F55871.D	05/16/06	22:19	NA	NA	NA	NA
12	P3886 AR 1248	F55872.D	05/16/06	22:34	NA	NA	NA	NA
13	P3994 AR 1254	F55873.D	05/16/06	22:48	NA	NA	NA	NA
14	P3887 AR 1262	F55874.D	05/16/06	23:03	NA	NA	NA	NA
15	P3888 AR 1268	F55875.D	05/16/06	23:17	NA	NA	NA	NA
16	P4057 AR 1254 0.0	F55876.D	05/16/06	23:47	NA	NA	NA	NA
17	P4056 AR 1254 0.0	F55877.D	05/16/06	0:01	NA	NA	NA	NA
18	P4055 AR 1254 0.1	F55878.D	05/17/06	0:16	NA	NA	NA	NA
19	P4054 AR 1254 0.2	F55879.D	05/17/06	0:31	NA	NA	NA	NA
20	P4053 AR 1254 0.5	F55880.D	05/17/06	0:45	NA	NA	NA	NA
21	P4052 AR 1254 1.0	F55881.D	05/17/06	0:59	NA	NA	NA	NA
22	P4010 AR 1254 0.5	F55882.D	05/17/06	1:13	2.64	10.37	2.72	10.02
23	ZZZZZ	F55883.D	5/17/06	1:28	2.64	10.37	2.72	10.02
24	ZZZZZ	F55885.D	5/17/06	2:12	2.64	10.37	2.72	10.02
25	ZZZZZ	F55886.D	5/17/06	2:26	2.64	10.37	2.72	10.02
26	ZZZZZ	F55895.D	5/17/06	4:53	2.64	10.38	2.73	10.03
27	ZZZZZ	F55896.D	5/17/06	5:07	2.64	10.39	2.73	10.03
28	ZZZZZ	F55897.D	5/17/06	5:22	2.64	10.39	2.73	10.03
29	P4021 16/60 0.5	F55900.D	5/17/06	6:06	2.64	10.39	2.72	10.04
30	P4053 AR 1254 0.5	F55901.D	5/17/06	6:21	NA	NA	NA	NA
31								

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PCB SYSTEM MONITORING COMPOUND RETENTION TIME SUMMARY

Quote#: SDG: 56617

Project No: 218736 Site ID: Location: ME

GC Columns: RTX-CLPesticidesI ID: 0.32(mm)

Instrument ID: F RTX-CLPesticidesII

Date Analyzed: 5/24/06

Lab File ID (Standard) : F56057 Time Analyzed: 15:30

					COLUMN #1		COLUMN #1	
					SMC1	SMC 2	SMC1	SMC 2
	PESTICIDE STANDARD				2.64	10.42	2.72	10.06
	UPPER LIMIT				2.69	10.47	2.77	10.11
	LOWER LIMIT				2.59	10.37	2.67	10.01
					COLUMN #1		COLUMN #2	
	SAMPLE ID	LAB FILE ID	DATE	TIME	SMC1 #	SMC 2 #	SMC1 #	SMC 2 #
01	P4020 16/60 0.25	F56057.D	05/24/06	15:30	2.64	10.42	2.72	10.06
02	B05246PAS, A/C	F56058.D	05/24/06	17:13	2.69	10.46	2.77	10.10
03	L05246PASB, A/C	F56059.D	05/24/06	17:28	2.65	10.43	2.73	10.07
04	LD05246PASB, A/C	F56060.D	05/24/06	17:42	2.65	10.42	2.73	10.06
05	ZZZZZ	F56061.D	05/24/06	17:57	2.64	10.41	2.73	10.06
06	ZZZZZ	F56062.D	05/24/06	18:11	2.64	10.41	2.72	10.06
07	ZZZZZ	F56063.D	05/24/06	18:26	2.64	10.40	2.72	10.05
08	ZZZZZ	F56064.D	05/24/06	18:41	2.64	10.41	2.72	10.05
09	ZZZZZ	F56065.D	05/24/06	19:10	2.64	10.40	2.72	10.05
10	ZZZZZ	F56066.D	05/24/06	19:25	2.64	10.40	2.72	10.05
11	56606-6, MS, A/C	F56067.D	05/24/06	19:39	2.64	10.40	2.72	10.05
12	56606-6, MSD, A/C	F56068.D	05/24/06	19:54	2.63	10.40	2.72	10.05
13	P4021 16/60 0.5	F56069.D	05/24/06	20:08	2.64	10.41	2.72	10.05
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PCB SYSTEM MONITORING COMPOUND RETENTION TIME SUMMARY

Quote#: SDG: 56617

Project No: 218736 Site ID: Location: ME

GC Columns: RTX-CLPesticidesI ID: 0.32(mm)

Instrument ID: F RTX-CLPesticidesII

Date Analyzed: 5/25/06

Lab File ID (Standard) : F56078 Time Analyzed: 14:26

					COLUMN #1		COLUMN #1	
					SMC1	SMC 2	SMC1	SMC 2
	PESTICIDE STANDARD				2.65	10.42	2.73	10.07
	UPPER LIMIT				2.70	10.47	2.78	10.12
	LOWER LIMIT				2.60	10.37	2.68	10.02
	SAMPLE ID	LAB FILE ID	DATE	TIME	COLUMN #1 SMC1 #	COLUMN #1 SMC 2 #	COLUMN #2 SMC1 #	COLUMN #2 SMC 2 #
01	P4020 16/60 0.25	F56078.D	05/25/06	14:26	2.65	10.42	2.73	10.07
02	B05256PWIPE, A/C	F56079.D	05/25/06	15:10	2.69	10.47	2.77	10.11
03	L05256PWIPE, A/C	F56080.D	05/25/06	15:25	2.65	10.44	2.73	10.08
04	LD05256PWIPE, A/	F56081.D	05/25/06	15:39	2.65	10.43	2.73	10.07
05	56617-24, A/C	F56082.D	05/25/06	16:08	2.69	10.47	2.76	10.11
06	56617-25, A/C	F56083.D	05/25/06	16:23	2.65	10.44	2.74	10.08
07	56617-26, A/C	F56084.D	05/25/06	16:52	2.65	10.43	2.73	10.08
08	56617-27, A/C	F56085.D	05/25/06	17:07	2.65	10.43	2.73	10.07
09	ZZZZZ	F56086.D	05/25/06	17:22	2.65	10.42	2.73	10.07
10	56617-29, A/C	F56087.D	05/25/06	17:36	2.65	10.42	2.73	10.07
11	56617-30, A/C	F56088.D	05/25/06	17:51	2.64	10.41	2.73	10.06
12	56617-31, A/C	F56089.D	05/25/06	18:05	2.64	10.40	2.73	10.05
13	56617-32, A/C	F56090.D	05/25/06	18:20	2.64	10.40	2.73	10.04
14	56617-33, A/C	F56091.D	05/25/06	18:35	2.64	10.40	2.72	10.04
15	56617-34, A/C	F56092.D	05/25/06	18:49	2.64	10.40	2.72	10.05
16	56617-35, A/C	F56093.D	05/25/06	19:19	2.64	10.40	2.72	10.05
17	56617-36, A/C	F56094.D	05/25/06	19:33	2.64	10.40	2.72	10.05
18	56617-37, A/C	F56095.D	05/25/06	19:48	2.64	10.41	2.73	10.06
19	P4021 16/60 0.5	F56096.D	05/25/06	20:03	2.64	10.41	2.72	10.05
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analytics environmental laboratory LLC

PCB SYSTEM MONITORING COMPOUND RETENTION TIME SUMMARY

Quote#: SDG: 56617

Project No: 218736 Site ID: Location: ME

GC Columns: RTX-CLPesticidesI ID: 0.32(mm)

Instrument ID: F RTX-CLPesticidesII

Date Analyzed: 5/25/06

Lab File ID (Standard) : F56096 Time Analyzed: 20:03

					COLUMN #1		COLUMN #1	
					SMC1	SMC 2	SMC1	SMC 2
	PESTICIDE STANDARD				2.64	10.41	2.72	10.05
	UPPER LIMIT				2.69	10.46	2.77	10.10
	LOWER LIMIT				2.59	10.36	2.67	10.00
	SAMPLE ID	LAB FILE ID	DATE	TIME	COLUMN #1 SMC1 #	COLUMN #1 SMC 2 #	COLUMN #2 SMC1 #	COLUMN #2 SMC 2 #
01	P4021 16/60 0.5	F56096.D	05/25/06	20:03	2.64	10.41	2.72	10.05
02	B05256PAS, A/C	F56097.D	05/25/06	20:17	2.64	10.41	2.73	10.06
03	L05256PAS, A/C	F56098.D	05/25/06	20:32	2.64	10.42	2.73	10.06
04	56617-1, A/C	F56099.D	05/25/06	20:46	2.64	10.41	2.73	10.06
05	56617-2, A/C	F56100.D	05/25/06	21:01	2.64	10.42	2.73	10.07
06	56617-6, A/C	F56101.D	05/25/06	21:16	2.64	10.42	2.73	10.06
07	56617-7, A/C	F56102.D	05/25/06	21:45	2.64	10.41	2.73	10.06
08	56617-8, A/C	F56103.D	05/25/06	22:00	2.64	10.41	2.72	10.06
09	56617-9, A/C	F56104.D	05/25/06	22:14	2.64	10.40	2.72	10.05
10	56617-10, A/C	F56105.D	05/25/06	22:29	2.64	10.40	2.72	10.04
11	56617-11, A/C	F56106.D	05/25/06	22:44	2.64	10.39	2.72	10.04
12	56617-12, A/C	F56107.D	05/25/06	22:58	2.64	10.39	2.72	10.04
13	56617-13, A/C	F56108.D	05/25/06	23:13	2.64	10.39	2.72	10.03
14	P4021 16/60 0.5	F56109.D	05/25/06	23:27	2.64	10.38	2.72	10.03
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analytics environmental laboratory LLC

PCB SYSTEM MONITORING COMPOUND RETENTION TIME SUMMARY

Quote#: SDG: 56617

Project No: 218736 Site ID: Location: ME

GC Columns: RTX-CLPesticidesI ID: 0.32(mm)

Instrument ID: F RTX-CLPesticidesII

Date Analyzed: 5/25/06

Lab File ID (Standard) : F56109 Time Analyzed: 23:27

					COLUMN #1		COLUMN #1	
					SMC1	SMC 2	SMC1	SMC 2
	PESTICIDE STANDARD				2.64	10.38	2.72	10.03
	UPPER LIMIT				2.69	10.43	2.77	10.08
	LOWER LIMIT				2.59	10.33	2.67	9.98
	SAMPLE ID	LAB FILE ID	DATE	TIME	COLUMN #1 SMC1 #	COLUMN #1 SMC 2 #	COLUMN #2 SMC1 #	COLUMN #2 SMC 2 #
01	P4021 16/60 0.5	F56109.D	05/25/06	23:27	2.64	10.38	2.72	10.03
02	ZZZZZ	F56110.D	05/25/06	23:42	2.64	10.39	2.72	10.04
03	ZZZZZ	F56111.D	05/25/06	23:57	2.64	10.38	2.72	10.03
04	56617-4, 1:10, A/C	F56112.D	05/26/06	0:26	2.64	10.38	2.72	10.03
05	56617-5, A/C	F56113.D	05/26/06	0:41	2.64	10.38	2.72	10.03
06	P4021 16/60 0.5	F56114.D	05/26/06	0:55	2.63	10.38	2.72	10.03
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PCB SYSTEM MONITORING COMPOUND RETENTION TIME SUMMARY

Quote#: SDG: 56617

Project No: 218736 Site ID: Location: ME

GC Columns: RTX-CLPesticidesI ID: 0.32(mm)

Instrument ID: F RTX-CLPesticidesII

Date Analyzed: 5/26/06

Lab File ID (Standard) : F56114 Time Analyzed: 0:55

				COLUMN #1		COLUMN #1	
				SMC1	SMC 2	SMC1	SMC 2
	PESTICIDE STANDARD			2.63	10.38	2.72	10.03
	UPPER LIMIT			2.68	10.43	2.77	10.08
	LOWER LIMIT			2.58	10.33	2.67	9.98
	SAMPLE ID	LAB FILE ID	DATE	TIME	COLUMN #1 SMC1 #	COLUMN #1 SMC 2 #	COLUMN #2 SMC1 #	COLUMN #2 SMC 2 #
01	P4021 16/60 0.5	F56114.D	05/26/06	0:55	2.63	10.38	2.72	10.03
02	ZZZZZ	F56115.D	05/26/06	7:01	2.70 *	10.43	2.70	10.09 *
03	B05266PAS, A/C	F56116.D	05/26/06	7:30	2.64	10.42	2.72	10.06
04	L05266PAS, A/C	F56117.D	05/26/06	7:45	2.67	10.43	2.74	10.08
05	LD05266PAS, A/C	F56118.D	05/26/06	8:00	2.64	10.42	2.73	10.07
06	56617-15, A/C	F56119.D	05/26/06	8:18	2.64	10.42	2.73	10.07
07	ZZZZZ	F56120.D	05/26/06	8:32	2.64	10.41	2.73	10.06
08	ZZZZZ	F56121.D	05/26/06	8:47	2.64	10.42	2.73	10.06
09	56617-18, A/C	F56122.D	05/26/06	9:01	2.64	10.42	2.72	10.07
10	56617-19, A/C	F56123.D	05/26/06	9:31	2.67	10.44 *	2.75	10.08
11	56617-20, A/C	F56124.D	05/26/06	9:45	2.64	10.42	2.73	10.07
12	56617-21, A/C	F56125.D	05/26/06	10:00	2.64	10.42	2.73	10.07
13	56617-22, A/C	F56126.D	05/26/06	10:27	2.68	10.46 *	2.76	10.09 *
14	56617-23, A/C	F56127.D	05/26/06	10:42	2.65	10.44 *	2.73	10.08
15	56617-23, MS, A/C	F56128.D	05/26/06	10:57	2.65	10.43	2.74	10.07
16	56617-23, MSD, A/C	F56129.D	05/26/06	11:11	2.64	10.43	2.73	10.07
17	P4021 16/60 0.5	F56130.D	05/26/06	11:41	2.64	10.43	2.73	10.08
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analytics environmental laboratory LLC

PCB SYSTEM MONITORING COMPOUND RETENTION TIME SUMMARY

Quote#: SDG: 56617

Project No: 218736 Site ID: Location: ME

GC Columns: RTX-CLPesticidesI ID: 0.32(mm)

Instrument ID: F RTX-CLPesticidesII

Date Analyzed: 5/26/06

Lab File ID (Standard) : F56132 Time Analyzed: 13:38

					COLUMN #1		COLUMN #1	
					SMC1	SMC 2	SMC1	SMC 2
	PESTICIDE STANDARD				2.66	10.45	2.74	10.09
	UPPER LIMIT				2.71	10.50	2.79	10.14
	LOWER LIMIT				2.61	10.40	2.69	10.04
	SAMPLE ID	LAB FILE ID	DATE	TIME	COLUMN #1 SMC1 #	COLUMN #1 SMC 2 #	COLUMN #2 SMC1 #	COLUMN #2 SMC 2 #
01	ZZZZZ	F56131.D	05/26/06	13:23	2.64	10.43	2.73	10.08
02	P4020 16/60 0.25	F56132.D	05/26/06	13:38	2.66	10.45	2.74	10.09
03	B05266PAS,RR A/C	F56133.D	05/26/06	14:07	2.68	10.47	2.76	10.11
04	56617-28, 1:10, A/C	F56134.D	05/26/06	14:22	2.65	10.45	2.73	10.09
05	56617-14, 1:2000, A	F56135.D	05/26/06	14:36	2.65	10.37 *	2.73	10.08
06	56617-3, 1:5, A/C	F56136.D	05/26/06	14:51	2.64	10.43	2.73	10.08
07	56617-16, 1:10, A/C	F56137.D	05/26/06	15:05	2.65	10.43	2.73	10.07
08	56617-17, 1:50, A/C	F56138.D	05/26/06	15:35	2.64	10.43	2.72	10.08
09	P4021 16/60 0.5	F56139.D	05/26/06	15:49	2.67	10.46	2.75	10.10
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Chain Of Custody Form

analytics environmental laboratory LLC

195 Commerce Way Suite E
Portsmouth, NH 03801
Phone (603) 436-5111
Fax (603) 430-2151

For Analytics Use Only Rev. 2, 10/25/04

Samples were:
1) Shipped or Hand-delivered
2) Temp blank °C 2.0
3) Received in good condition Y or N
4) pH checked by: NA
5) Labels checked by: EAM 5/25/00

Project#: 218736 Proj. Name: GP
Company: W&C
Contact: Karl Kasper
Address: 41 Hutchins DR.
Phone: 774-2112 PO# Quote #
Sampler (Signature): Josh Bowe

Matrix Key:
WW=Wastewater
SW=Surfacewater
GW=Groundwater
DW=Drinkingwater
S=Soil/Sludge
O=Oil
E=Extract
X=Other

Container Key
P=plastic G=glass

Preservation

Station Identification	Sample Date	Sample Time	Analysis	Unpres	4°C	HNO3	H2SO4	HCL	Methanol	Other	Matrix	Container number/type		pH	Analytics Sample #
SB-2 (0-4)	5/24	1620	PCB		X										56617-1
SB-3 (0-4)	"	1630	"		X										-2
SB-4 (0-4)	"	1635	"		X										-3
SB-10 (0-4)	"	1040	"		X										-4
SB-11 (0-4)	"	1040	"		X										-5
SB-12 (0-4)	"	1140	"		X										-6
SB-15 (0-4)	"	1400	"		X										-7
SB-18 (0-4)	"	16:00	"		X										-8
SB-25 (0-4)	"	13:15	"		X										-9
SB-26 (0-4)	"	15:01	"		X										-10
SB-28 (0-4)	"	1545	"		X										-11
															-12

MTO 6/25/00

Email RESULTS? YES NO
Email _____
Fax RESULTS? YES NO
Fax# _____

Comments / Instructions:

Turnaround Request
Standard ☐ Due Date _____
Priority ☐ Due Date _____

Page 1 of 3

Relinquished By Sampler:	Relinquished By:	Relinquished By:
Josh Bowe		
Date: 5-24-00	Date: 5-25	Date:
Time: 17:19	Time: 11:15 am	Time:
Received By:	Received By:	Received By:

Analytics\AEL Documents\AEL COC

226/228

Chain Of Custody Form

analytics environmental laboratory LLC

195 Commerce Way Suite E
Portsmouth, NH 03801
Phone (603) 436-5111
Fax (603) 430-2151

For Analytics Use Only Rev. 2, 10/25/04

Samples were:
1) Shipped or hand-delivered
2) Temp blank °C 2.0
3) Received in good condition Y or N
4) pH checked by: NA
5) Labels checked by: DPM 5/25/06

Project#: Proj. Name:
Company:
Contact:
Address:
Phone: PO# Quote #
Sampler (Signature): [signature]

Matrix Key:
WW=Wastewater
SW=Surfacewater
GW=Groundwater
DW=Drinkingwater
S=Soil/Sludge
O=Oil
F=Extract
X=Other

Container Key
P=plastic G=glass

Preservation

Station Identification	Sample Date	Sample Time	Analysis	Unpres	4°C	HNO_3	H_2SO_4	HCL	Methanol	Other	Matrix	Container number/type	pH	Analytics Sample #
SC-01	5/24	1020	PCB		X									56617-12
SC-02	"	1025	"		X									-13
SC-03	"	1432	"		X									-14
SS-05	"	0920	"		X									-15
SS-06	"	0930	"		X									-16
SS-07	"	0935	"		X									-17
SS-08	"	0915	"		X									-18
SS-09	"	0905	"		X									-19
SS-09 DUP	"	0905	"		X									-20
SS-10	"	0855	"		X									-21
SS-11	"	0850	"		X									-22
SS-12	"	0845	"		X									-23

Email RESULTS? YES NO
Email
Fax RESULTS? YES NO
Fax#

Comments / Instructions:

Turnaround Request
Standard ☐ Due Date
Priority ☐ Due Date

Page 2 of 3

Received By: [signature] Time: 17:19 Date: 5/24/06
Received By: [signature] Time: 11:15am Date: 5/25
Received By: Time: Date:
Relinquished By Sampler: Josh Bowe
Relinquished By: [signature]
Relinquished By:

Analytics\AEL Documents\AEL COC

227/228

Katahdin ANALYTICAL SERVICES
340 County Road No. 5
P.O. Box 720
Westbrook, ME 04092
Tel: (207) 874-2400
Fax: (207) 775-4029

CHAIN of CUSTODY

PLEASE BEAR DOWN AND PRINT LEGIBLY IN PEN

Client: WTC	Contact: Karl Kasp	Phone #: ( ) 774-2112	Fax #: ( )
Address: 41 Hutchins Dr	City: Portland	State: ME	Zip Code: 04102
Purchase Order #: 218736	Proj. Name / No.: GP	Katahdin Quote #	
Bill (if different than above)	Address		
Sampler (Print / Sign)		Copies To:	

LAB USE ONLY | WORK ORDER #:

KATAHDIN PROJECT NUMBER

REMARKS:

SHIPPING INFO: ☐ FED EX ☐ UPS ☐ CLIENT

AIRBILL NO:

TEMP°C ☐ TEMP BLANK ☐ INTACT ☐ NOT INTACT

*	Sample Description	Date / Time coll'd	Matrix	No. of Cntrs.	ANALYSIS AND CONTAINER TYPE PRESERVATIVES: PCB (Filt. ☐Y ☐N)									
	W-1	5/24 / 1122	Wipe	1	✓							56617-24		
	W-2	" / 1120	"	"	✓									-25
	W-3	" / 1130	"	"	✓									-26
	W-4	" / 1115	"	"	✓									-27
	W-5	" / 1105	"	"	✓									-28
	W-6	" / 1100	"	"	✓									-29
	W-6Dup	" / 1100	"	"	✓									-30
	W-7	" / 1058	"	"	✓									-31
	W-8	" / 1055	"	"	✓									-32
	W-9	" / 1050	"	"	✓									-33
	W-10	" / 1045	"	"	✓									-34
	W-11	" / 1032	"	"	✓									-35
	W-12	" / 1037	"	"	✓									-36
	W-13	" / 1945	"	"	✓									-37
		/												
		/												

COMMENTS

emails DPM 5/25/06

Received at AEL temp 2.0°C MJG 5.25.06

Relinquished By: (Signature)	Date / Time	Received By: (Signature)	Relinquished By: (Signature)	Date / Time	Received By: (Signature)
Josh Bowe	5/24 17:18	E Kirkhurst	E Kirkhurst	5/25 11:15	
Relinquished By: (Signature)	Date / Time	Received By: (Signature)	Relinquished By: (Signature)	Date / Time	Received By: (Signature)

THE TERMS AND CONDITIONS ON THE REVERSE SIDE HEREOF SHALL GOVERN SERVICES EXCEPT WHEN A SIGNED CONTRACTUAL AGREEMENT EXISTS

analytics environmental laboratory LLC

195 Commerce Way Suite E
Portsmouth, New Hampshire 03801
603-436-5111 Fax 603-430-2151
800-929-9906
www.analyticslab.com

June 6, 2006

Mr. Karl Kasper
Woodard & Curran, Inc.
41 Hutchins drive
Portland ME 04102

RE: **Analytical Results Case Narrative**
Analytics # 56637

Dear Mr. Kasper;

Enclosed please find the analytical results for samples submitted for the above-mentioned project. The attached Cover Page lists the sample IDs, Lab tracking numbers and collection dates for the samples included in this deliverable.

Samples were analyzed VOCs by EPA Method 8260B, SVOCs by EPA Method 8270C, Polychlorinated Biphenyls (PCBs) by EPA Method 8082, and RCRA 8 Metals. Metals analyses were subcontracted to Maine Environmental Laboratory (MEL)

Unless otherwise noted in the Non-conformance Summary listed below, all of the quality control (QC) criteria including initial calibration, calibration verification, surrogate recovery, holding time and method accuracy/precision for these analyses were within acceptable limits.

This Level III data package has been assembled in the following order:

Case Narrative/Non-Conformance Summary
Sample Log Sheet - Cover Page
VOC Form I Data Sheet for Samples and Blanks
VOC Form 2 Surrogate Recoveries
VOC Form 3 MS/MSD (LCS) Recoveries
VOC Form 4 Method Blank Summary (equiv.) GC/MS Logbook Sheets
VOC Form 5 DFTPP Tune Summary
VOC Form 6 Initial Calibration Data
VOC Form 7 Continuing Calibration Check
VOC Form 8 Internal Standard Area and RT Summary
SVOC Form I Data Sheet for Samples and Blanks
SVOC Form 2 Surrogate Recoveries
SVOC Form 3 MS/MSD (LCS) Recoveries
SVOC Form 4 Method Blank Summary (equiv.) GC/MS Logbook Sheets
SVOC Form 5 DFTPP Tune Summary
SVOC Form 6 Initial Calibration Data
SVOC Form 7 Continuing Calibration Check
SVOC Form 8 Internal Standard Area and RT Summary

PCB Form 1 Data Sheet for Samples and Blanks
PCB Form 10 Confirmation Results Form (if required)
PCB Form 2 Surrogate Recoveries
PCB Form 3 MS/MSD (LCS) Recoveries
PCB Form 4 Method Blank Summary (equiv.) GC/MS Logbook Sheets
PCB Form 6 Initial Calibration Data
PCB Form 7 Continuing Calibration Check
PCB Form 8 System Monitoring Compound RT Summary
Metals Data Package from Maine Environmental
Chain of Custody (COC) Forms and Sample Receipt Checklist

QC NON CONFORMANCE SUMMARY

Sample Receipt:
No QC deviations.

EPA Method 8260B Volatile Organics:
The following analytes were not "J" flagged in this report: Methylene chloride, Diethyl ether, Acetone, Hexachlorobutadiene, and Naphthalene.

Seven analytes had a %RSD's greater than 15% for the initial calibration analyzed 05/22/06 (V805226B). The average RSD of all analytes was <15%. These analytes all had acceptable recoveries in the second source calibration check standard. Results were reported without qualification.

The laboratory control sample (L805306B) analyzed 05/30/06 had two analytes (Vinyl Chloride and 1,2-Dibromo-3-chloropropane) outside the laboratory acceptance criteria. The MS/MSD analyzed on 56637-2 (MW-4) was in control for all analytes, however the RPD for Bromomethane had high recovery. None of the analytes with QC deviations were detected in the samples associated with this QC and results were reported without qualification.

No extra sample was provided for MS/MSD analysis for solid matrix. Laboratory control samples were used to assess accuracy and precision.

Sample 56637-14 (SB-31 (0-4)) had surrogate recoveries below the laboratory acceptance criteria. The sample was reanalyzed with similar results. The original analysis is reported in this report with a comment to this affect.

Two analytes Bromomethane and Bromoform) had a %RSD's greater than 15% for the initial calibration analyzed 05/26/06 (V805266C). The average RSD of all analytes was <15%. These analytes all had acceptable recoveries in the second source calibration check standard. Results were reported without qualification.

2,2-Dichloropropane had %D above the acceptance criteria in the continuing calibration standard (file C59118SC) analyzed 05/30/06. Vinyl Chloride had low recovery in the laboratory control sample (LS05306C). The laboratory control sample was reanalyzed (LS05306C2) with similar results for vinyl chloride. Three other analytes also had slightly low recovery. (see Form III) Two analytes had high RPD (2,2-Dichloropropane and Chloroform). These analytes were not detected in any samples for this SDG and results were reported without qualification.

Semivolatile Organic Compounds:
The following analytes were not 'J" flagged in this report: Benzidine, Bis(2-ethyl hexyl) phthalate, Diethyl phthalate, Dimethyl phtahlate, Di-n-butyl phthalate, Di-n-octyl phthalate and Butyl benzyl phthalate.

Benzidine had high recovery in the second source check standard. The quantitation limit for Benzidine is estimated due to the QC criteria not being met. A comment to this affect was placed on the data summary pages.

No extra sample was provided for MS/MSD analysis for aqueous matrix. Laboratory control samples were used to assess accuracy and precision.

Three aqueous samples had some low surrogate recoveries. (See Form II) All the samples were re-injected and reanalyzed to confirm the results. The original analysis is reported in this report with a comment to this affect.

Several solid samples had surrogate recovery failures, internal standard failures or calibration range exceedances on the original analysis. All of the samples were reanalyzed and the results of both analysis are contained in this report. The original results were flagged accordingly and contain comments to this affect.

The aqueous laboratory control sample duplicate (LD05306AW) had six analytes with recovery below the laboratory acceptance criteria. The RPDs for these six compounds were above the laboratory acceptance criteria. The laboratory control sample duplicate was reinjected with similar results. Results were reported without qualification except as stated above.

Several analytes had recoveries outside the laboratory acceptance criteria in the MS/MSD analyzed on 56637-23 (Pile 2). The laboratory control samples (L 05306AS/LD05306AS) were in control for all analytes. Results were reported without qualification except as stated above.

PCBs by EPA Method 8082:
Sample 56637-6 (SB-21 (4-8')) had % difference between the analytical columns above the laboratory acceptance criteria. Results were reported with a "D" flag and a comment to this affect.

The laboratory blank (B05306PAS) had retention times for of the surrogates slightly outside the established windows. Results were reported without qualification.

Sincerely,
ANALYTICS Environmental Laboratory, LLC

Stephen Knollmeyer
Laboratory Director

analytics environmental laboratory LLC

195 Commerce Way Suite E
Portsmouth, New Hampshire 03801
603-436-5111 Fax 603-430-2151
800-929-9906
www.analyticslab.com

Mr. Karl Kasper
Woodard & Curran
41 Hutchins Drive
Portland ME 04102

Report Number: 56637

Revision: Rev. 0

Re: GP Old Town Mill **218736.002**

Enclosed are the results of the analyses on your sample(s). Samples were received on 27 May 2006 and analyzed for the tests listed below. Samples were received in acceptable condition, with the exceptions noted below or on the chain of custody. The results reported herein conform to the most current NELAC standards, where applicable, unless otherwise narrated in the body of the report. Please see individual reports for specific methodologies and references.

Lab Number	Sample Date	Station Location	Analysis	Comments
56637-1	05/26/06	MW-5	EPA 8260 Volatile Organics	
	05/26/06	MW-5	EPA 8270 Acid/Base Neutrals	
	05/26/06	MW-5	RCRA Metals	
56637-2	05/26/06	MW-4	EPA 8260 Volatile Organics	
	05/26/06	MW-4	EPA 8270 Acid/Base Neutrals	
	05/26/06	MW-4	RCRA Metals	
56637-3	05/26/06	SB-37	EPA 8260 Volatile Organics	
	05/26/06	SB-37	EPA 8270 Acid/Base Neutrals	
	05/26/06	SB-37	RCRA Metals	
56637-4	05/25/06	SB-22 (0-4)	EPA 8082 (PCBs only)	
	05/25/06	SB-22 (0-4)	EPA 8260 Volatile Organics	
	05/25/06	SB-22 (0-4)	EPA 8270 Acid/Base Neutrals	
	05/25/06	SB-22 (0-4)	RCRA Metals	
56637-5	05/25/06	SB-1 (6-10')	EPA 8260 Volatile Organics	
56637-6	05/25/06	SB-21 (4-8')	EPA 8082 (PCBs only)	

Sample Receipt Exceptions: None

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If you have any further question on the analytical methods or these results, do not hesitate to call.

Authorized signature _____
Stephen L. Knollmeyer Lab. Director

Date 5/31/2006

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analytics environmental laboratory LLC

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800-929-9906
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Mr. Karl Kasper
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Lab Number	Sample Date	Station Location	Analysis	Comments
56637-7	05/25/06	SB-35 (0-6")	EPA 8082 (PCBs only)	
56637-8	05/25/06	SB-5 (4-6')	EPA 8082 (PCBs only)	
56637-9	05/25/06	SB-21 (0-4')	EPA 8082 (PCBs only)	
	05/25/06	SB-21 (0-4')	EPA 8260 Volatile Organics	
	05/25/06	SB-21 (0-4')	EPA 8270 Acid/Base Neutrals	
	05/25/06	SB-21 (0-4')	RCRA Metals	
56637-10	05/25/06	SB-5 (0-4')	EPA 8082 (PCBs only)	
	05/25/06	SB-5 (0-4')	EPA 8260 Volatile Organics	
	05/25/06	SB-5 (0-4')	EPA 8270 Acid/Base Neutrals	
	05/25/06	SB-5 (0-4')	RCRA Metals	
56637-11	05/25/06	SB-6 (0-4')	EPA 8260 Volatile Organics	
	05/25/06	SB-6 (0-4')	EPA 8270 Acid/Base Neutrals	
	05/25/06	SB-6 (0-4')	RCRA Metals	
56637-12	05/25/06	SB-23 (0-4')	EPA 8082 (PCBs only)	
	05/25/06	SB-23 (0-4')	EPA 8270 Acid/Base Neutrals	

Sample Receipt Exceptions: None

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Lab Number	Sample Date	Station Location	Analysis	Comments
	05/25/06	SB-23 (0-4')	RCRA Metals	
56637-13	05/25/06	SB-24 (0-4')	EPA 8082 (PCBs only)	
56637-14	05/25/06	SB-31 (0-4)	EPA 8260 Volatile Organics	
	05/25/06	SB-31 (0-4)	EPA 8270 Acid/Base Neutrals	
	05/25/06	SB-31 (0-4)	RCRA Metals	
56637-15	05/25/06	SB-30 (0-4)	EPA 8260 Volatile Organics	
	05/25/06	SB-30 (0-4)	EPA 8270 Acid/Base Neutrals	
	05/25/06	SB-30 (0-4)	RCRA Metals	
56637-16	05/25/06	SB-38 (0-4)	EPA 8260 Volatile Organics	
	05/25/06	SB-38 (0-4)	EPA 8270 Acid/Base Neutrals	
	05/25/06	SB-38 (0-4)	RCRA Metals	
56637-17	05/25/06	SB-37 (0-4)	EPA 8260 Volatile Organics	
	05/25/06	SB-37 (0-4)	EPA 8270 Acid/Base Neutrals	
	05/25/06	SB-37 (0-4)	RCRA Metals	
56637-18	05/25/06	SB-15 (0-4)	EPA 8260 Volatile Organics	

Sample Receipt Exceptions: None

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Lab Number	Sample Date	Station Location	Analysis	Comments
	05/25/06	SB-15 (0-4)	EPA 8270 Acid/Base Neutrals	
	05/25/06	SB-15 (0-4)	RCRA Metals	
56637-19	05/26/06	SB-41 (0-4)	EPA 8082 (PCBs only)	
	05/26/06	SB-41 (0-4)	EPA 8260 Volatile Organics	
	05/26/06	SB-41 (0-4)	EPA 8270 Acid/Base Neutrals	
	05/26/06	SB-41 (0-4)	RCRA Metals	
56637-20	05/26/06	SB-15	EPA 8260 Volatile Organics	
	05/26/06	SB-15	EPA 8270 Acid/Base Neutrals	
	05/26/06	SB-15	RCRA Metals	
56637-21	05/26/06	MW-3	EPA 8260 Volatile Organics	
	05/26/06	MW-3	EPA 8270 Acid/Base Neutrals	
	05/26/06	MW-3	RCRA Metals	
56637-22	05/26/06	Pile 1	EPA 8270 Acid/Base Neutrals	
	05/26/06	Pile 1	RCRA Metals	
56637-23	05/26/06	Pile 2	EPA 8270 Acid/Base Neutrals	

Sample Receipt Exceptions: None

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Lab Number	Sample Date	Station Location	Analysis	Comments
	05/26/06	Pile 2	RCRA Metals	
56637-24	05/26/06	SB-40 (1 1/2-5 1/2)	EPA 8260 Volatile Organics	
56637-25	05/25/06	SB-38 DUP	EPA 8260 Volatile Organics	
56637-26	05/26/06	MW-5	EPA 8260 Volatile Organics	
56637-27	05/25/06	SB-30	EPA 8260 Volatile Organics	
56637-28	05/25/06	SB-15	EPA 8260 Volatile Organics	
56637-29	05/26/06	Trip Blank	Electronic Data Deliverable	
	05/26/06	Trip Blank	EPA 8260 Volatile Organics	

Sample Receipt Exceptions: None

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Date 5/31/2006

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VOLATILE
DATA SUMMARIES

analytics environmental laboratory LLC

195 Commerce Way
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800-929-9906

Mr. Karl Kasper
Woodard & Curran
41 Hutchins Drive
Portland ME 04102

June 6, 2006

CLIENT SAMPLE ID

Project Name: GP Old Town Mill

Project Number:

Field Sample ID: LAB QC

SAMPLE DATA

Lab Sample ID: B805306B
Matrix: Aqueous
Percent Solid: N/A
Dilution Factor: 1
Collection Date: N/A
Lab Receipt Date: N/A
Analysis Date: 05/30/06

ANALYTICAL RESULTS VOLATILE ORGANICS

COMPOUND	Quantitation Limit µg/L	Result µg/L	COMPOUND	Quantitation Limit µg/L	Result µg/L
Benzene	2	U	1,3-Dichloropropane	2	U
Bromobenzene	2	U	cis-1,3-Dichloropropene	2	U
Bromochloromethane	2	U	trans-1,3-Dichloropropene	2	U
Bromodichloromethane	2	U	2,2-Dichloropropane	2	U
Bromoform	2	U	1,1-Dichloropropene	2	U
Bromomethane	2	U	Ethylbenzene	2	U
n-butylbenzene	2	U	Hexachlorobutadiene	2	U
sec-butylbenzene	2	U	Isopropylbenzene	2	U
tert-butylbenzene	2	U	p-isopropyltoluene	2	U
Carbon Tetrachloride	2	U	Methylene Chloride	5	U
Chlorobenzene	2	U	Methyl-tert-butyl ether (MTBE)	2	U
Chloroethane	2	U	Naphthalene	2	U
Chloroform	2	U	n-Propylbenzene	2	U
Chloromethane	2	U	Styrene	2	U
2-Chlorotoluene	2	U	1,1,1,2-Tetrachloroethane	2	U
4-Chlorotoluene	2	U	1,1,2,2-Tetrachloroethane	2	U
Dibromochloromethane	2	U	Tetrachloroethene	2	U
1,2-Dibromo-3-chloropropane	2	U	Toluene	2	U
1,2-Dibromoethane	2	U	1,2,3-Trichlorobenzene	2	U
Dibromomethane	2	U	1,2,4-Trichlorobenzene	2	U
1,2-Dichlorobenzene	2	U	1,1,1-Trichloroethane	2	U
1,3-Dichlorobenzene	2	U	1,1,2-Trichloroethane	2	U
1,4-Dichlorobenzene	2	U	Trichloroethene	2	U
Dichlorodifluoromethane	2	U	Trichlorofluoromethane	2	U
1,1-Dichloroethane	2	U	1,2,3-Trichloropropane	2	U
1,2-Dichloroethane	2	U	1,2,4-Trimethylbenzene	2	U
1,1-Dichloroethene	2	U	1,3,5-Trimethylbenzene	2	U
cis-1,2-Dichloroethene	2	U	Vinyl Chloride	2	U
trans-1,2-Dichloroethene	2	U	o-Xylene	2	U
1,2-Dichloropropane	2	U	m,p-Xylene	2	U
Acetone	10	U	Diethyl ether	2	U
Carbon Disulfide	2	U	2-Hexanone	10	U
Tetrahydrofuran	5	U	Methyl isobutyl ketone	10	U
Methyl ethyl ketone	10	U	Di-isopropyl ether (DIPE)	2	U
t-Butyl alcohol (TBA)	20	U	Ethyl t-butyl ether (ETBE)	2	U
t-Amyl methyl ether (TAME)	2	U			

Surrogate Standard Recovery

d4-1,2-Dichloroethane	99 %	d8-Toluene	99 %	Bromofluorobenzene	100 %

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis was conducted according to: Test Methods for Evaluating Solid Waste, SW-846 Method 8260B.

COMMENTS:

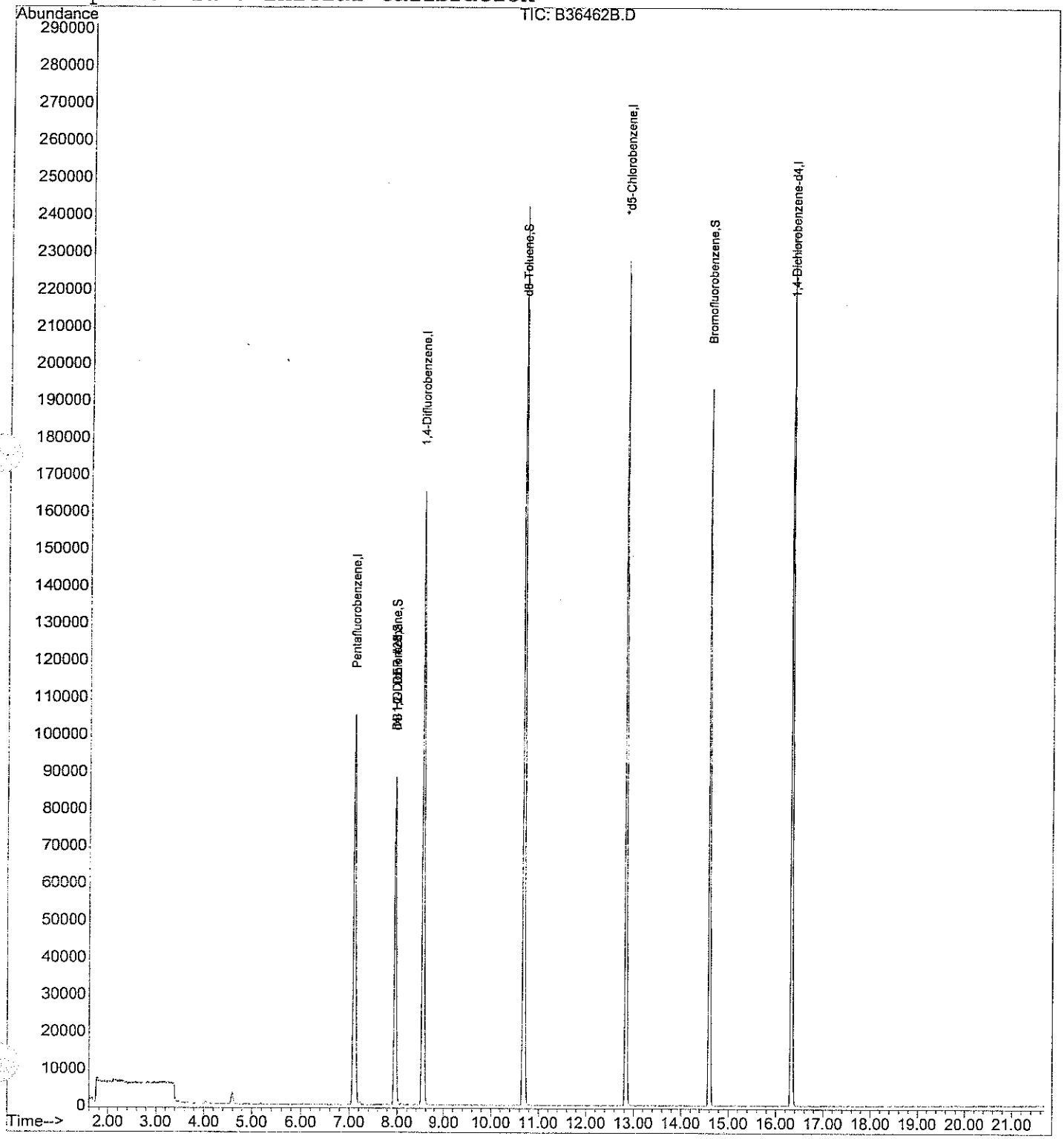
8260 full

Authorized signature

Quantitation Report

Data File : C:\HPCHEM\1\DATA\053006-B\B36462B.D Vial: 6
Acq On : 30 May 2006 10:51 am Operator:
Sample : B805306B Inst : Instrumen
Misc : 5000 Multiplr: 1.00
MS Integration Params: rteint.p
Quant Time: May 30 11:12 2006 Quant Results File: V805266B.RES

Method : C:\HPCHEM\1\METHODS\V805266B.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Tue May 30 08:22:32 2006
Response via : Initial Calibration



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Mr. Karl Kasper
Woodard & Curran
41 Hutchins Drive
Portland ME 04102

June 6, 2006

CLIENT SAMPLE ID

Project Name: GP Old Town Mill

Project Number: 218736.002

Field Sample ID: LAB QC

SAMPLE DATA

Lab Sample ID: MB05306C
Matrix: Solid
Percent Solid: 100
Dilution Factor: 100
Collection Date: N/A
Lab Receipt Date: N/A
Analysis Date: 05/30/06

ANALYTICAL RESULTS VOLATILE ORGANICS

COMPOUND	Quantitation Limit µg/kg	Result µg/kg	COMPOUND	Quantitation Limit µg/kg	Result µg/kg
Benzene	100	U	1,3-Dichloropropane	100	U
Bromobenzene	100	U	cis-1,3-Dichloropropene	100	U
Bromochloromethane	100	U	trans-1,3-Dichloropropene	100	U
Bromodichloromethane	75	U	2,2-Dichloropropane	100	U
Bromoform	75	U	1,1-Dichloropropene	100	U
Bromomethane	100	U	Ethylbenzene	100	U
n-butylbenzene	100	U	Hexachlorobutadiene	100	U
sec-butylbenzene	100	U	Isopropylbenzene	100	U
tert-butylbenzene	100	U	p-isopropyltoluene	100	U
Carbon Tetrachloride	100	U	Methylene Chloride	500	U
Chlorobenzene	100	U	Methyl-tert-butyl ether (MTBE)	100	U
Chloroethane	100	U	Naphthalene	100	U
Chloroform	75	U	n-Propylbenzene	100	U
Chloromethane	100	U	Styrene	100	U
2-Chlorotoluene	100	U	1,1,1,2-Tetrachloroethane	100	U
4-Chlorotoluene	100	U	1,1,2,2-Tetrachloroethane	75	U
Dibromochloromethane	75	U	Tetrachloroethene	100	U
1,2-Dibromo-3-chloropropane	100	U	Toluene	100	U
1,2-Dibromoethane	75	U	1,2,3-Trichlorobenzene	100	U
Dibromomethane	100	U	1,2,4-Trichlorobenzene	100	U
1,2-Dichlorobenzene	100	U	1,1,1-Trichloroethane	100	U
1,3-Dichlorobenzene	100	U	1,1,2-Trichloroethane	75	U
1,4-Dichlorobenzene	100	U	Trichloroethene	100	U
Dichlorodifluoromethane	100	U	Trichlorofluoromethane	100	U
1,1-Dichloroethane	100	U	1,2,3-Trichloropropane	100	U
1,2-Dichloroethane	75	U	1,2,4-Trimethylbenzene	100	U
1,1-Dichloroethene	75	U	1,3,5-Trimethylbenzene	100	U
cis-1,2-Dichloroethene	100	U	Vinyl Chloride	100	U
trans-1,2-Dichloroethene	100	U	o-Xylene	100	U
1,2-Dichloropropane	75	U	m,p-Xylene	100	U
Acetone	1000	U	Diethyl ether	100	U
Carbon Disulfide	100	U	2-Hexanone	1000	U
Tetrahydrofuran	500	U	Methyl isobutyl ketone	1000	U
Methyl ethyl ketone	1000	U	Di-isopropyl ether (DIPE)	100	U
t-Butyl alcohol (TBA)	2000	U	Ethyl t-butyl ether (ETBE)	100	U
t-Amyl methyl ether (TAME)	100	U			

Surrogate Standard Recovery

d4-1,2-Dichloroethane	84	%	d8-Toluene	87	%	Bromofluorobenzene	90 %

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis was conducted according to: Test Methods for Evaluating Solid Waste, SW-846 Method 8260B.

COMMENTS: Results are expressed on a dry weight basis.

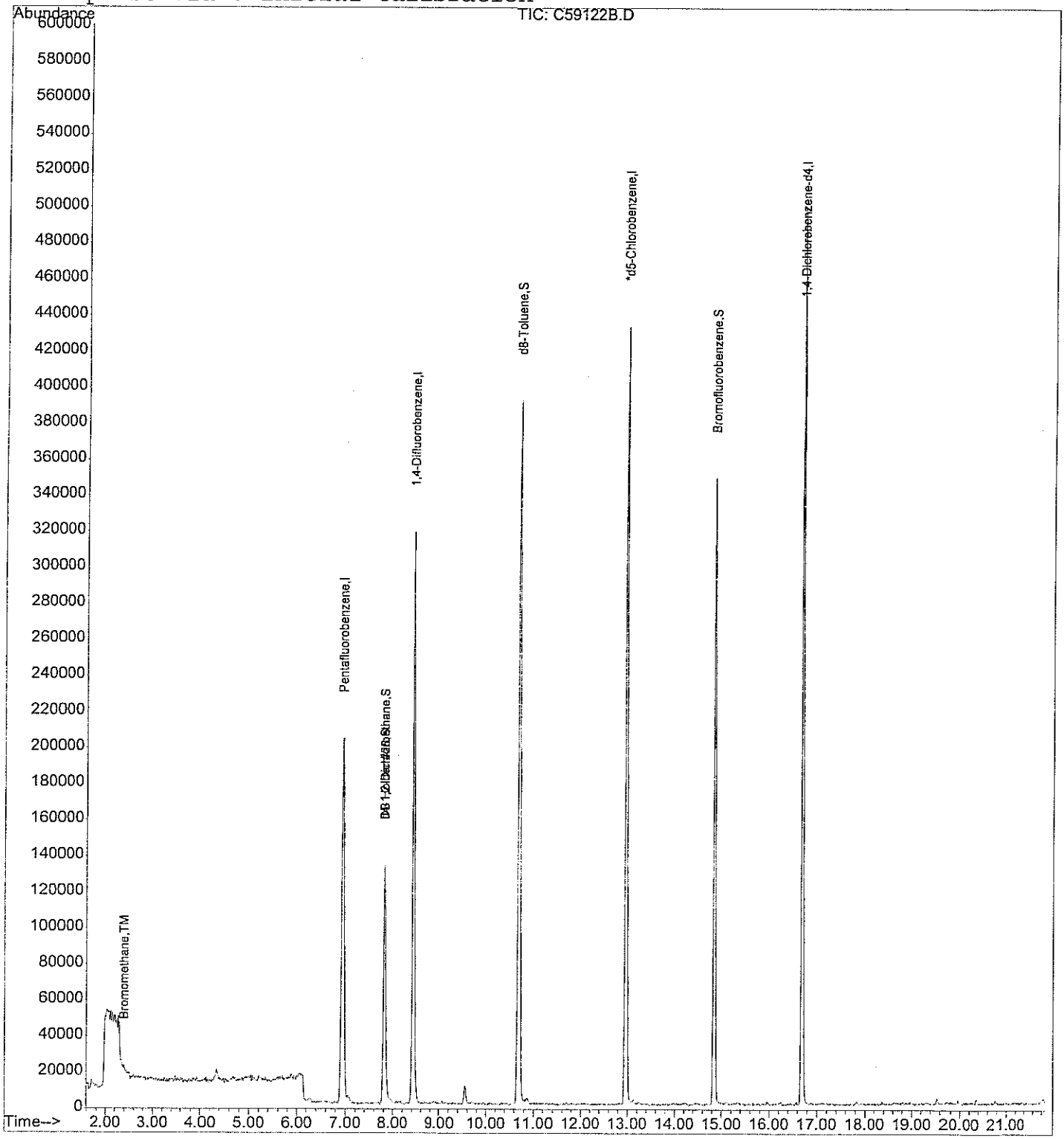
8260 full

Authorized signature

Quantitation Report

Data File : D:\HPCHEM\1\DATA\053006C\C59122B.D Vial: 7
Acq On : 30 May 2006 10:55 am Operator:
Sample : MB05306C Inst : Instr_C
Misc : 50,5.00,SOIL Multiplr: 1.00
MS Integration Params: rteint.p
Quant Time: May 30 10:59 2006 Quant Results File: V805226C.RES

Method : D:\HPCHEM\1\METHODS\V805226C.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Mon May 22 14:18:09 2006
Response via : Initial Calibration



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Mr. Karl Kasper
Woodard & Curran
41 Hutchins Drive
Portland ME 04102

May 31, 2006

CLIENT SAMPLE ID

Project Name: GP Old Town Mill

Project Number: 218736.002

Field Sample ID: MW-5

SAMPLE DATA

Lab Sample ID: 56637-1
Matrix: Aqueous
Percent Solid: N/A
Dilution Factor: 1
Collection Date: 05/26/06
Lab Receipt Date: 05/27/06
Analysis Date: 05/30/06

ANALYTICAL RESULTS VOLATILE ORGANICS

COMPOUND	Quantitation Limit µg/L	Result µg/L	COMPOUND	Quantitation Limit µg/L	Result µg/L
Benzene	2	U	1,3-Dichloropropane	2	U
Bromobenzene	2	U	cis-1,3-Dichloropropene	2	U
Bromochloromethane	2	U	trans-1,3-Dichloropropene	2	U
Bromodichloromethane	2	U	2,2-Dichloropropane	2	U
Bromoform	2	U	1,1-Dichloropropene	2	U
Bromomethane	2	U	Ethylbenzene	2	U
n-butylbenzene	2	U	Hexachlorobutadiene	2	U
sec-butylbenzene	2	U	Isopropylbenzene	2	U
tert-butylbenzene	2	U	p-isopropyltoluene	2	U
Carbon Tetrachloride	2	U	Methylene Chloride	5	U
Chlorobenzene	2	U	Methyl-tert-butyl ether (MTBE)	2	U
Chloroethane	2	U	Naphthalene	2	U
Chloroform	2	U	n-Propylbenzene	2	U
Chloromethane	2	U	Styrene	2	U
2-Chlorotoluene	2	U	1,1,1,2-Tetrachloroethane	2	U
4-Chlorotoluene	2	U	1,1,2,2-Tetrachloroethane	2	U
Dibromochloromethane	2	U	Tetrachloroethene	2	U
1,2-Dibromo-3-chloropropane	2	U	Toluene	2	U
1,2-Dibromoethane	2	U	1,2,3-Trichlorobenzene	2	U
Dibromomethane	2	U	1,2,4-Trichlorobenzene	2	U
1,2-Dichlorobenzene	2	U	1,1,1-Trichloroethane	2	U
1,3-Dichlorobenzene	2	U	1,1,2-Trichloroethane	2	U
1,4-Dichlorobenzene	2	U	Trichloroethene	2	U
Dichlorodifluoromethane	2	U	Trichlorofluoromethane	2	U
1,1-Dichloroethane	2	U	1,2,3-Trichloropropane	2	U
1,2-Dichloroethane	2	U	1,2,4-Trimethylbenzene	2	U
1,1-Dichloroethene	2	U	1,3,5-Trimethylbenzene	2	U
cis-1,2-Dichloroethene	2	U	Vinyl Chloride	2	U
trans-1,2-Dichloroethene	2	U	o-Xylene	2	U
1,2-Dichloropropane	2	U	m,p-Xylene	2	U
Acetone	10	U	Diethyl ether	2	U
Carbon Disulfide	2	U	2-Hexanone	10	U
Tetrahydrofuran	5	U	Methyl isobutyl ketone	10	U
Methyl ethyl ketone	10	U	Di-isopropyl ether (DIPE)	2	U
t-Butyl alcohol (TBA)	20	U	Ethyl t-butyl ether (ETBE)	2	U
t-Amyl methyl ether (TAME)	2	U			

Surrogate Standard Recovery

d4-1,2-Dichloroethane	101	%	d8-Toluene	100	%	Bromofluorobenzene	99 %

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis was conducted according to: Test Methods for Evaluating Solid Waste, SW-846 Method 8260B.

COMMENTS:

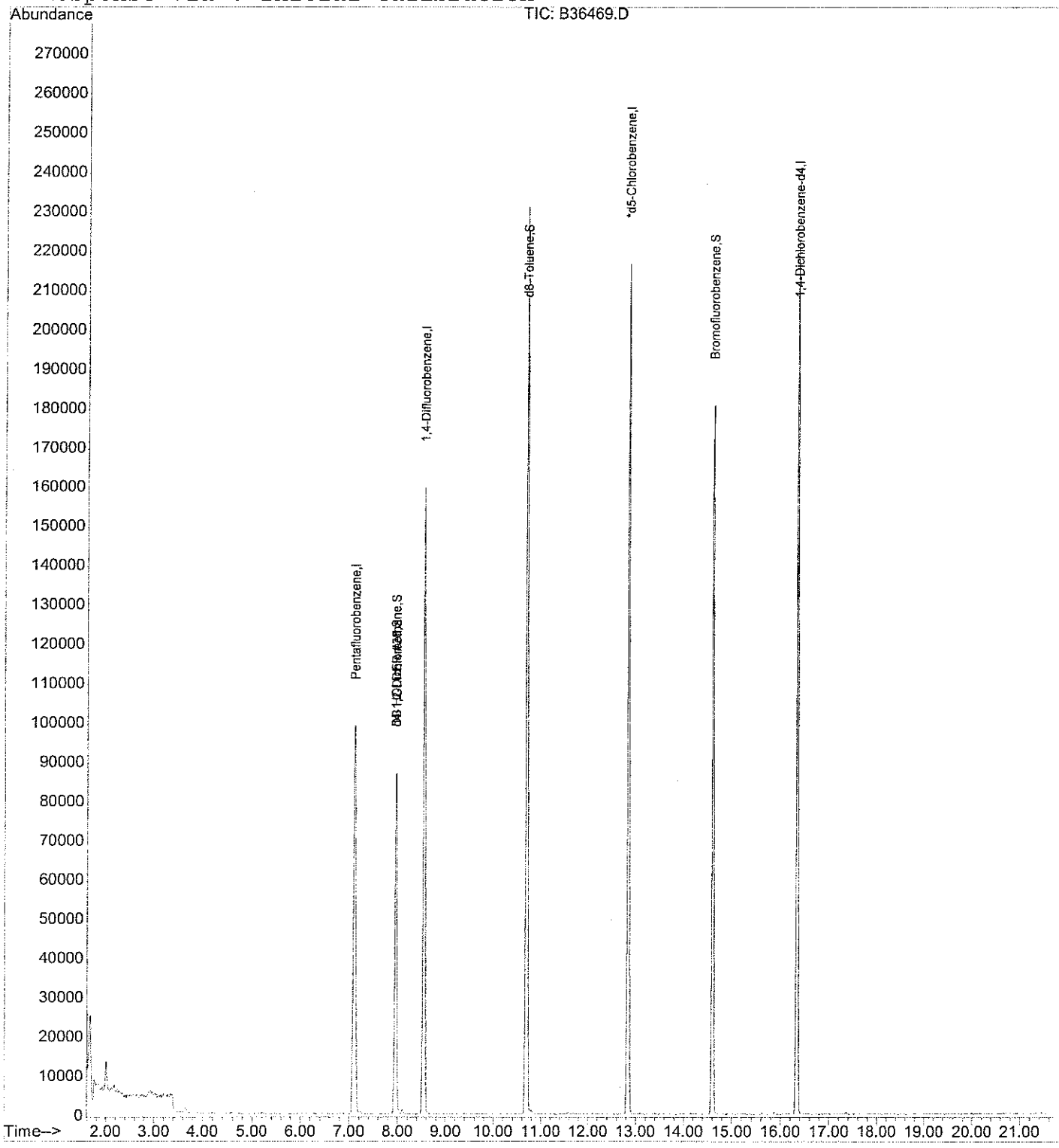
8260 full

Authorized signature

Quantitation Report

Data File : C:\HPCHEM\1\DATA\053006-B\B36469.D Vial: 13
Acq On : 30 May 2006 2:25 pm Operator:
Sample : 56637-1 Inst : Instrumen
Misc : 5000 Multiplr: 1.00
MS Integration Params: rteint.p
Quant Time: May 30 14:39 2006 Quant Results File: V805266B.RES

Method : C:\HPCHEM\1\METHODS\V805266B.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Tue May 30 08:22:32 2006
Response via : Initial Calibration



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Mr. Karl Kasper
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41 Hutchins Drive
Portland ME 04102

May 31, 2006

CLIENT SAMPLE ID

Project Name: GP Old Town Mill

Project Number: 218736.002

Field Sample ID: MW-4

SAMPLE DATA

Lab Sample ID: 56637-2
Matrix: Aqueous
Percent Solid: N/A
Dilution Factor: 1
Collection Date: 05/26/06
Lab Receipt Date: 05/27/06
Analysis Date: 05/30/06

ANALYTICAL RESULTS VOLATILE ORGANICS

COMPOUND	Quantitation Limit µg/L	Result µg/L	COMPOUND	Quantitation Limit µg/L	Result µg/L
Benzene	2	U	1,3-Dichloropropane	2	U
Bromobenzene	2	U	cis-1,3-Dichloropropene	2	U
Bromochloromethane	2	U	trans-1,3-Dichloropropene	2	U
Bromodichloromethane	2	U	2,2-Dichloropropane	2	U
Bromoform	2	U	1,1-Dichloropropene	2	U
Bromomethane	2	U	Ethylbenzene	2	U
n-butylbenzene	2	U	Hexachlorobutadiene	2	U
sec-butylbenzene	2	U	Isopropylbenzene	2	U
tert-butylbenzene	2	U	p-isopropyltoluene	2	2
Carbon Tetrachloride	2	U	Methylene Chloride	5	U
Chlorobenzene	2	U	Methyl-tert-butyl ether (MTBE)	2	U
Chloroethane	2	U	Naphthalene	2	U
Chloroform	2	U	n-Propylbenzene	2	U
Chloromethane	2	U	Styrene	2	U
2-Chlorotoluene	2	U	1,1,1,2-Tetrachloroethane	2	U
4-Chlorotoluene	2	U	1,1,2,2-Tetrachloroethane	2	U
Dibromochloromethane	2	U	Tetrachloroethene	2	U
1,2-Dibromo-3-chloropropane	2	U	Toluene	2	U
1,2-Dibromoethane	2	U	1,2,3-Trichlorobenzene	2	U
Dibromomethane	2	U	1,2,4-Trichlorobenzene	2	U
1,2-Dichlorobenzene	2	U	1,1,1-Trichloroethane	2	U
1,3-Dichlorobenzene	2	U	1,1,2-Trichloroethane	2	U
1,4-Dichlorobenzene	2	U	Trichloroethene	2	U
Dichlorodifluoromethane	2	U	Trichlorofluoromethane	2	U
1,1-Dichloroethane	2	U	1,2,3-Trichloropropane	2	U
1,2-Dichloroethane	2	U	1,2,4-Trimethylbenzene	2	U
1,1-Dichloroethene	2	U	1,3,5-Trimethylbenzene	2	U
cis-1,2-Dichloroethene	2	U	Vinyl Chloride	2	U
trans-1,2-Dichloroethene	2	U	o-Xylene	2	U
1,2-Dichloropropane	2	U	m,p-Xylene	2	U
Acetone	10	13	Diethyl ether	2	U
Carbon Disulfide	2	U	2-Hexanone	10	U
Tetrahydrofuran	5	U	Methyl isobutyl ketone	10	U
Methyl ethyl ketone	10	U	Di-isopropyl ether (DIPE)	2	U
t-Butyl alcohol (TBA)	20	U	Ethyl t-butyl ether (ETBE)	2	U
t-Amyl methyl ether (TAME)	2	U			

Surrogate Standard Recovery

d4-1,2-Dichloroethane	102	%	d8-Toluene	100	%	Bromofluorobenzene	102 %

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis was conducted according to: Test Methods for Evaluating Solid Waste, SW-846 Method 8260B.

COMMENTS:

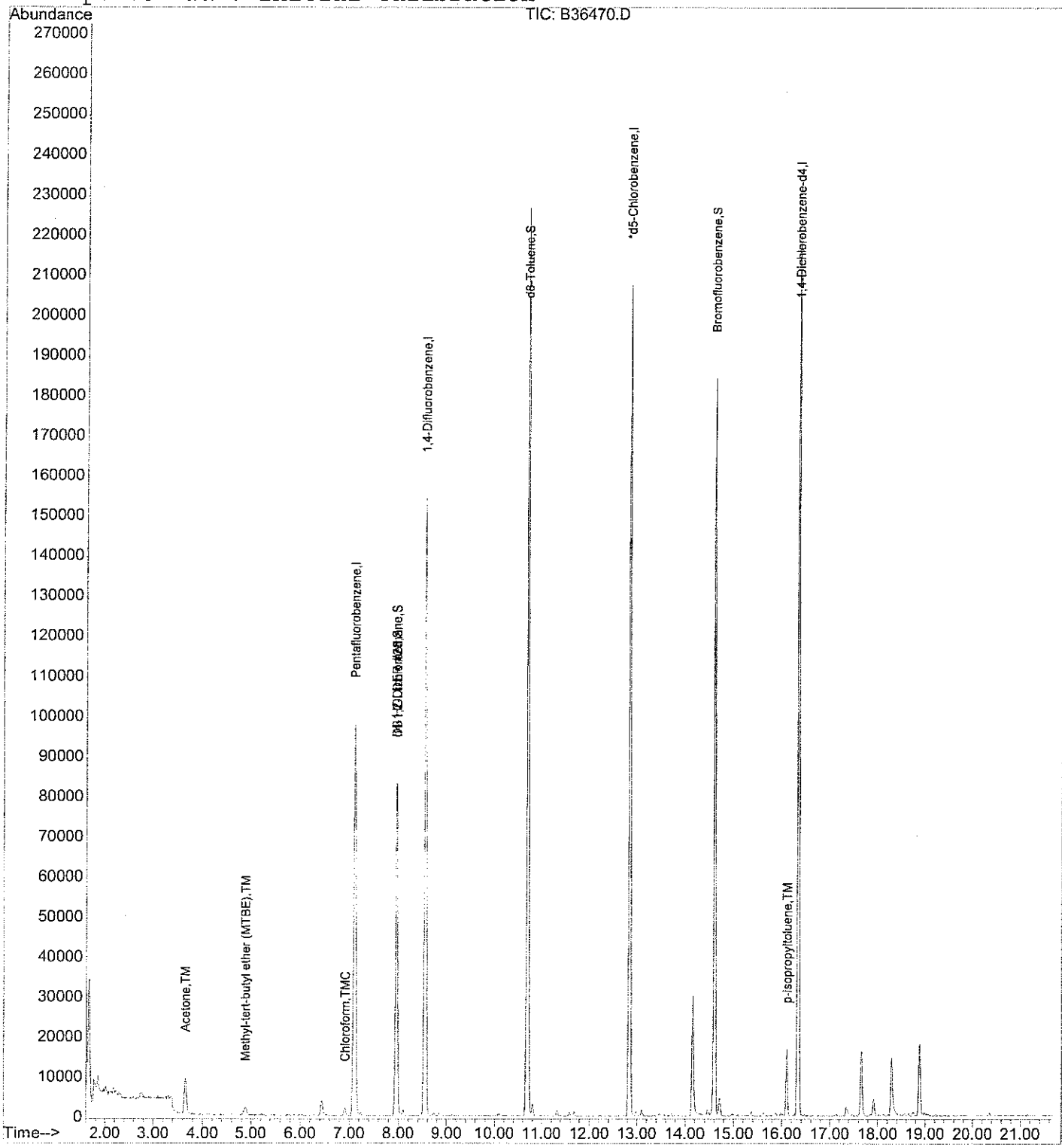
8260 full

Authorized signature

Quantitation Report

Data File : C:\HPCHEM\1\DATA\053006-B\B36470.D Vial: 14
Acq On : 30 May 2006 2:54 pm Operator:
Sample : 56637-2 Inst : Instrumen
Misc : 5000 Multiplr: 1.00
MS Integration Params: rteint.p
Quant Time: May 30 15:08 2006 Quant Results File: V805266B.RES

Method : C:\HPCHEM\1\METHODS\V805266B.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Tue May 30 08:22:32 2006
Response via : Initial Calibration



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May 31, 2006

CLIENT SAMPLE ID

Project Name: GP Old Town Mill

Project Number: 218736.002

Field Sample ID: SB-37

SAMPLE DATA

Lab Sample ID: 56637-3
Matrix: Aqueous
Percent Solid: N/A
Dilution Factor: 1
Collection Date: 05/26/06
Lab Receipt Date: 05/27/06
Analysis Date: 05/30/06

ANALYTICAL RESULTS VOLATILE ORGANICS

COMPOUND	Quantitation Limit µg/L	Result µg/L	COMPOUND	Quantitation Limit µg/L	Result µg/L
Benzene	2	U	1,3-Dichloropropane	2	U
Bromobenzene	2	U	cis-1,3-Dichloropropene	2	U
Bromochloromethane	2	U	trans-1,3-Dichloropropene	2	U
Bromodichloromethane	2	U	2,2-Dichloropropane	2	U
Bromoform	2	U	1,1-Dichloropropene	2	U
Bromomethane	2	U	Ethylbenzene	2	U
n-butylbenzene	2	U	Hexachlorobutadiene	2	U
sec-butylbenzene	2	U	Isopropylbenzene	2	U
tert-butylbenzene	2	U	p-isopropyltoluene	2	U
Carbon Tetrachloride	2	U	Methylene Chloride	5	U
Chlorobenzene	2	U	Methyl-tert-butyl ether (MTBE)	2	U
Chloroethane	2	U	Naphthalene	2	U
Chloroform	2	U	n-Propylbenzene	2	U
Chloromethane	2	U	Styrene	2	U
2-Chlorotoluene	2	U	1,1,1,2-Tetrachloroethane	2	U
4-Chlorotoluene	2	U	1,1,2,2-Tetrachloroethane	2	U
Dibromochloromethane	2	U	Tetrachloroethene	2	U
1,2-Dibromo-3-chloropropane	2	U	Toluene	2	U
1,2-Dibromoethane	2	U	1,2,3-Trichlorobenzene	2	U
Dibromomethane	2	U	1,2,4-Trichlorobenzene	2	U
1,2-Dichlorobenzene	2	U	1,1,1-Trichloroethane	2	U
1,3-Dichlorobenzene	2	U	1,1,2-Trichloroethane	2	U
1,4-Dichlorobenzene	2	U	Trichloroethene	2	U
Dichlorodifluoromethane	2	U	Trichlorofluoromethane	2	U
1,1-Dichloroethane	2	U	1,2,3-Trichloropropane	2	U
1,2-Dichloroethane	2	U	1,2,4-Trimethylbenzene	2	U
1,1-Dichloroethene	2	U	1,3,5-Trimethylbenzene	2	U
cis-1,2-Dichloroethene	2	U	Vinyl Chloride	2	U
trans-1,2-Dichloroethene	2	U	o-Xylene	2	U
1,2-Dichloropropane	2	U	m,p-Xylene	2	U
Acetone	10	U	Diethyl ether	2	U
Carbon Disulfide	2	U	2-Hexanone	10	U
Tetrahydrofuran	5	U	Methyl isobutyl ketone	10	U
Methyl ethyl ketone	10	U	Di-isopropyl ether (DIPE)	2	U
t-Butyl alcohol (TBA)	20	U	Ethyl t-butyl ether (ETBE)	2	U
t-Amyl methyl ether (TAME)	2	U			

Surrogate Standard Recovery

d4-1,2-Dichloroethane	102	%	d8-Toluene	99	%	Bromofluorobenzene	99 %

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis was conducted according to: Test Methods for Evaluating Solid Waste, SW-846 Method 8260B.

COMMENTS:

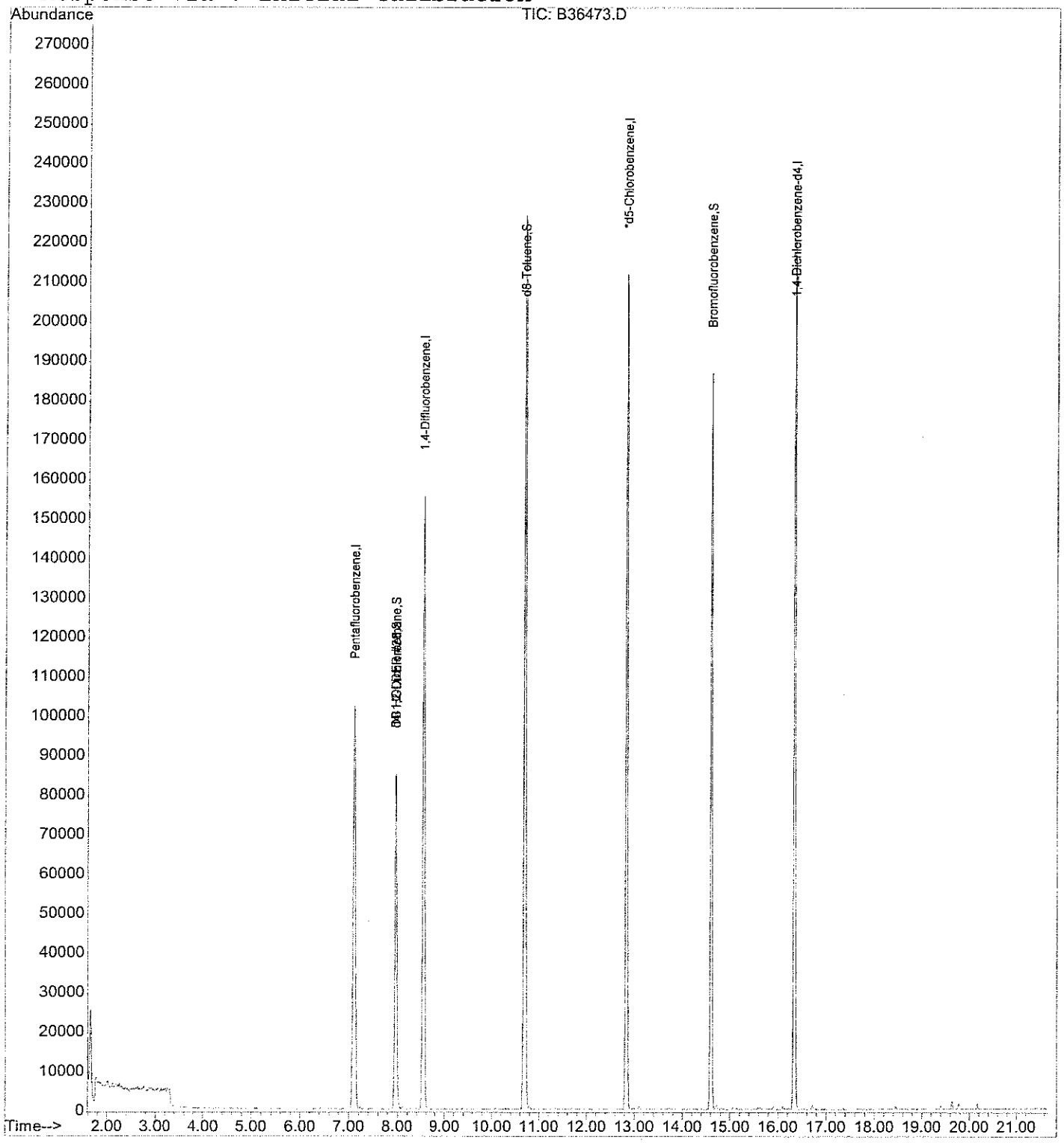
8260 full

Authorized signature

Quantitation Report

Data File : C:\HPCHEM\1\DATA\053006-B\B36473.D Vial: 17
Acq On : 30 May 2006 4:22 pm Operator:
Sample : 56637-3 Inst : Instrumen
Misc : 5000 Multiplr: 1.00
MS Integration Params: rteint.p
Quant Time: May 30 16:37 2006 Quant Results File: V805266B.RES

Method : C:\HPCHEM\1\METHODS\V805266B.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Tue May 30 08:22:32 2006
Response via : Initial Calibration



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May 31, 2006

CLIENT SAMPLE ID

Project Name: GP Old Town Mill

Project Number: 218736.002

Field Sample ID: SB-22 (0-4)

SAMPLE DATA

Lab Sample ID: 56637-4
Matrix: Solid
Percent Solid: 85
Dilution Factor: 135
Collection Date: 05/25/06
Lab Receipt Date: 05/27/06
Analysis Date: 05/30/06

ANALYTICAL RESULTS VOLATILE ORGANICS

COMPOUND	Quantitation Limit µg/kg	Result µg/kg	COMPOUND	Quantitation Limit µg/kg	Result µg/kg
Benzene	135	U	1,3-Dichloropropane	135	U
Bromobenzene	135	U	cis-1,3-Dichloropropene	135	U
Bromochloromethane	135	U	trans-1,3-Dichloropropene	135	U
Bromodichloromethane	101	U	2,2-Dichloropropane	135	U
Bromoform	101	U	1,1-Dichloropropene	135	U
Bromomethane	135	U	Ethylbenzene	135	U
n-butylbenzene	135	U	Hexachlorobutadiene	135	U
sec-butylbenzene	135	U	Isopropylbenzene	135	U
tert-butylbenzene	135	U	p-isopropyltoluene	135	U
Carbon Tetrachloride	135	U	Methylene Chloride	673	U
Chlorobenzene	135	U	Methyl-tert-butyl ether (MTBE)	135	U
Chloroethane	135	U	Naphthalene	135	U
Chloroform	101	U	n-Propylbenzene	135	U
Chloromethane	135	U	Styrene	135	U
2-Chlorotoluene	135	U	1,1,1,2-Tetrachloroethane	135	U
4-Chlorotoluene	135	U	1,1,2,2-Tetrachloroethane	101	U
Dibromochloromethane	101	U	Tetrachloroethene	135	U
1,2-Dibromo-3-chloropropane	135	U	Toluene	135	U
1,2-Dibromoethane	101	U	1,2,3-Trichlorobenzene	135	U
Dibromomethane	135	U	1,2,4-Trichlorobenzene	135	U
1,2-Dichlorobenzene	135	U	1,1,1-Trichloroethane	135	U
1,3-Dichlorobenzene	135	U	1,1,2-Trichloroethane	101	U
1,4-Dichlorobenzene	135	U	Trichloroethene	135	U
Dichlorodifluoromethane	135	U	Trichlorofluoromethane	135	U
1,1-Dichloroethane	135	U	1,2,3-Trichloropropane	135	U
1,2-Dichloroethane	101	U	1,2,4-Trimethylbenzene	135	U
1,1-Dichloroethene	101	U	1,3,5-Trimethylbenzene	135	U
cis-1,2-Dichloroethene	135	U	Vinyl Chloride	135	U
trans-1,2-Dichloroethene	135	U	o-Xylene	135	U
1,2-Dichloropropane	101	U	m,p-Xylene	135	U
Acetone	1350	U	Diethyl ether	135	U
Carbon Disulfide	135	U	2-Hexanone	1350	U
Tetrahydrofuran	673	U	Methyl isobutyl ketone	1350	U
Methyl ethyl ketone	1350	U	Di-isopropyl ether (DIPE)	135	U
t-Butyl alcohol (TBA)	2690	U	Ethyl t-butyl ether (ETBE)	135	U
t-Amyl methyl ether (TAME)	135	U			

Surrogate Standard Recovery

d4-1,2-Dichloroethane	78	%	d8-Toluene	76	%	Bromofluorobenzene	79 %

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis was conducted according to: Test Methods for Evaluating Solid Waste, SW-846 Method 8260B.

COMMENTS: Results are expressed on a dry weight basis. Sample collection and analysis in accordance with SW-846 method 5035A.

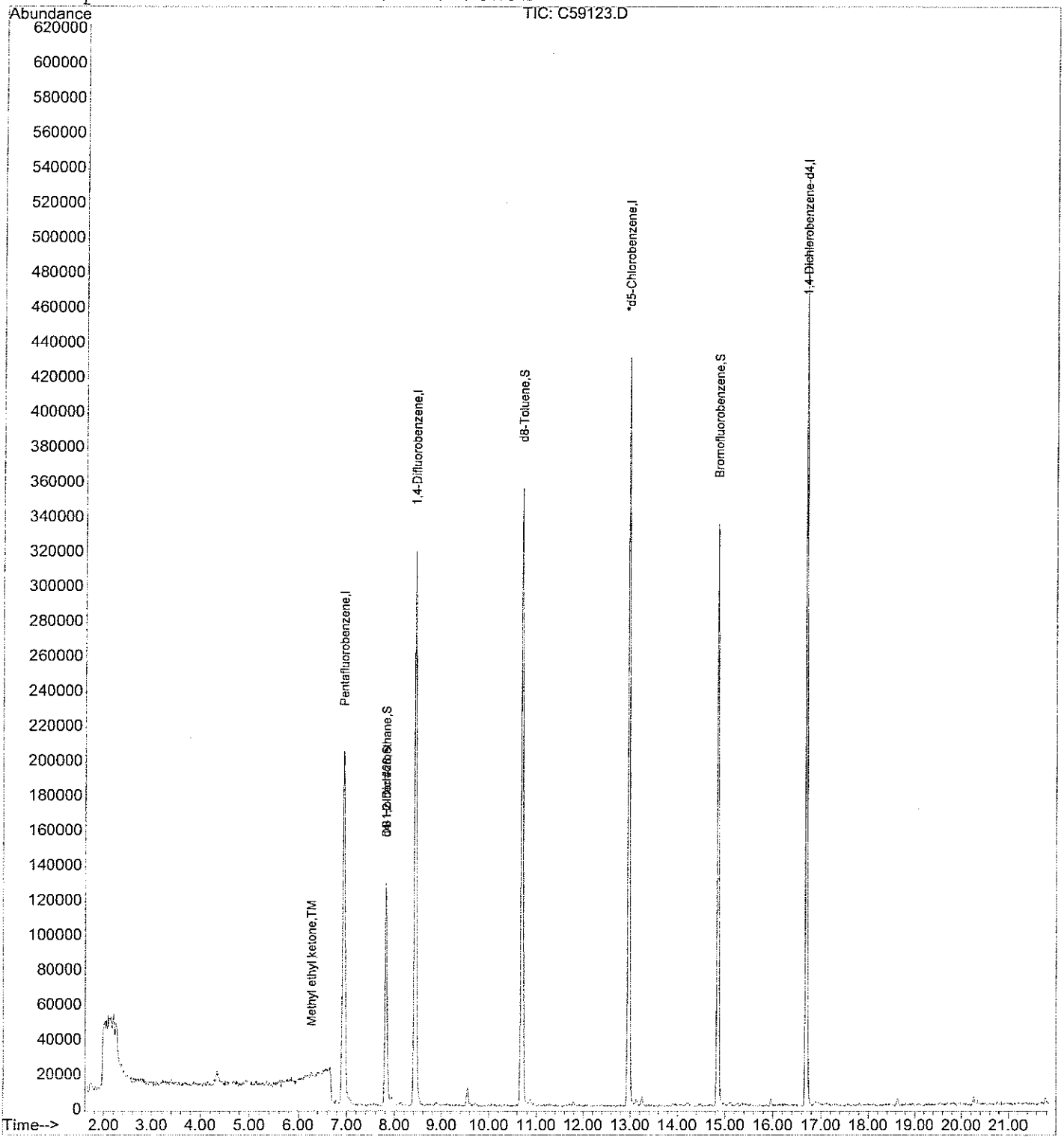
8260 full

Authorized signature

Quantitation Report

Data File : D:\HPCHEM\1\DATA\053006C\C59123.D Vial: 15
Acq On : 30 May 2006 11:30 am Operator:
Sample : 56637-4 Inst : Instr_C
Misc : 50,4.39,SOIL Multiplr: 1.00
MS Integration Params: rteint.p
Quant Time: May 30 11:36 2006 Quant Results File: V805226C.RES

Method : D:\HPCHEM\1\METHODS\V805226C.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Mon May 22 14:18:09 2006
Response via : Initial Calibration



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May 31, 2006

CLIENT SAMPLE ID

Project Name: GP Old Town Mill

Project Number: 218736.002

Field Sample ID: SB-1 (6-10')

SAMPLE DATA

Lab Sample ID: 56637-5
Matrix: Aqueous
Percent Solid: N/A
Dilution Factor: 1
Collection Date: 05/25/06
Lab Receipt Date: 05/27/06
Analysis Date: 05/30/06

ANALYTICAL RESULTS VOLATILE ORGANICS

COMPOUND	Quantitation Limit µg/L	Result µg/L	COMPOUND	Quantitation Limit µg/L	Result µg/L
Benzene	2	U	1,3-Dichloropropane	2	U
Bromobenzene	2	U	cis-1,3-Dichloropropene	2	U
Bromochloromethane	2	U	trans-1,3-Dichloropropene	2	U
Bromodichloromethane	2	U	2,2-Dichloropropane	2	U
Bromoform	2	U	1,1-Dichloropropene	2	U
Bromomethane	2	U	Ethylbenzene	2	U
n-butylbenzene	2	U	Hexachlorobutadiene	2	U
sec-butylbenzene	2	U	Isopropylbenzene	2	U
tert-butylbenzene	2	U	p-isopropyltoluene	2	U
Carbon Tetrachloride	2	U	Methylene Chloride	5	U
Chlorobenzene	2	U	Methyl-tert-butyl ether (MTBE)	2	U
Chloroethane	2	U	Naphthalene	2	U
Chloroform	2	U	n-Propylbenzene	2	U
Chloromethane	2	U	Styrene	2	U
2-Chlorotoluene	2	U	1,1,1,2-Tetrachloroethane	2	U
4-Chlorotoluene	2	U	1,1,2,2-Tetrachloroethane	2	U
Dibromochloromethane	2	U	Tetrachloroethene	2	U
1,2-Dibromo-3-chloropropane	2	U	Toluene	2	U
1,2-Dibromoethane	2	U	1,2,3-Trichlorobenzene	2	U
Dibromomethane	2	U	1,2,4-Trichlorobenzene	2	U
1,2-Dichlorobenzene	2	U	1,1,1-Trichloroethane	2	U
1,3-Dichlorobenzene	2	U	1,1,2-Trichloroethane	2	U
1,4-Dichlorobenzene	2	U	Trichloroethene	2	U
Dichlorodifluoromethane	2	U	Trichlorofluoromethane	2	U
1,1-Dichloroethane	2	U	1,2,3-Trichloropropane	2	U
1,2-Dichloroethane	2	U	1,2,4-Trimethylbenzene	2	U
1,1-Dichloroethene	2	U	1,3,5-Trimethylbenzene	2	U
cis-1,2-Dichloroethene	2	U	Vinyl Chloride	2	U
trans-1,2-Dichloroethene	2	U	o-Xylene	2	U
1,2-Dichloropropane	2	U	m,p-Xylene	2	U
Acetone	10	U	Diethyl ether	2	U
Carbon Disulfide	2	U	2-Hexanone	10	U
Tetrahydrofuran	5	U	Methyl isobutyl ketone	10	U
Methyl ethyl ketone	10	U	Di-isopropyl ether (DIPE)	2	U
t-Butyl alcohol (TBA)	20	U	Ethyl t-butyl ether (ETBE)	2	U
t-Amyl methyl ether (TAME)	2	U			

Surrogate Standard Recovery

d4-1,2-Dichloroethane	100	%	d8-Toluene	100	%	Bromofluorobenzene	100	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis was conducted according to: Test Methods for Evaluating Solid Waste, SW-846 Method 8260B.

COMMENTS:

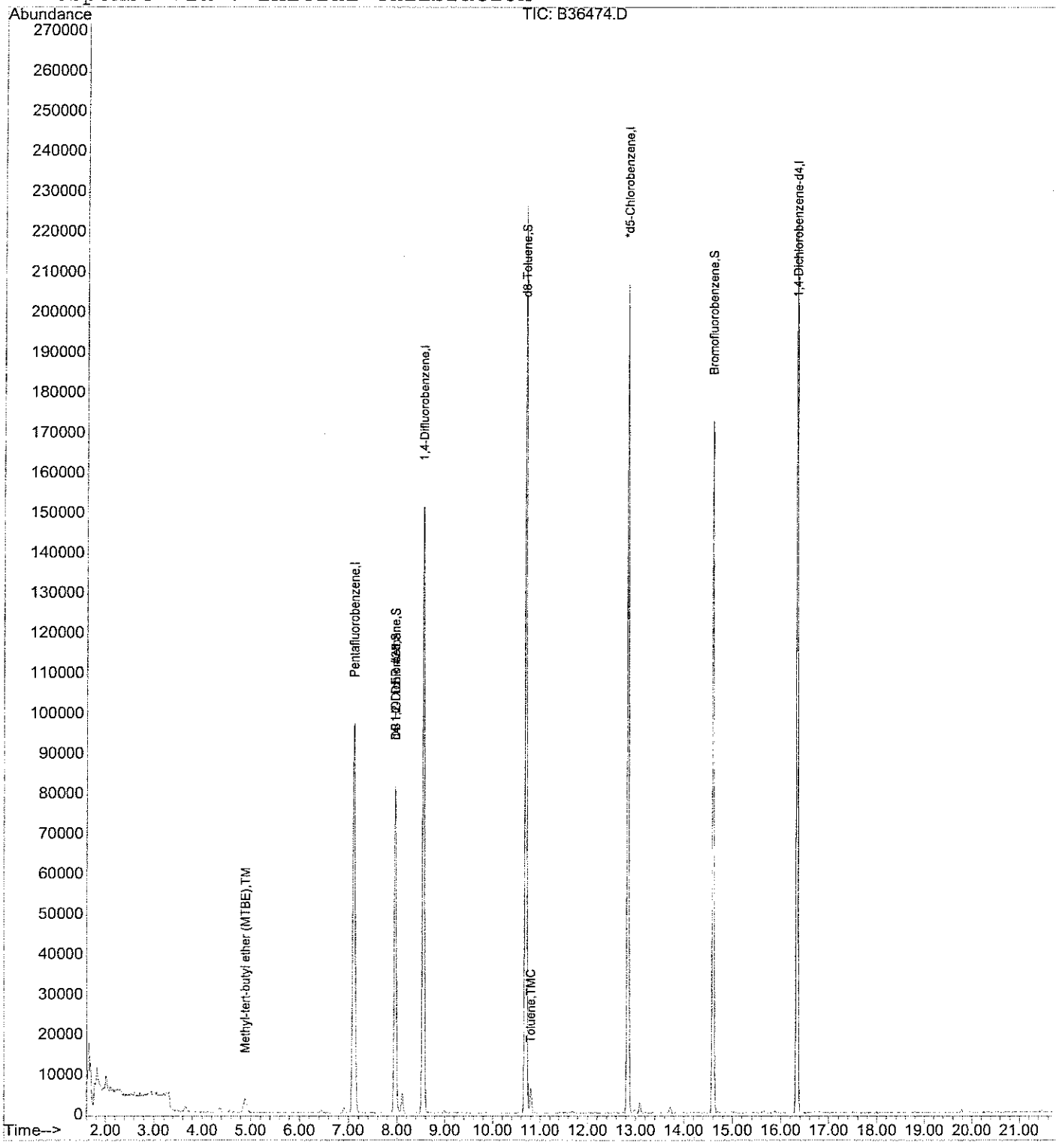
8260 full

Authorized signature

Quantitation Report

Data File : C:\HPCHEM\1\DATA\053006-B\B36474.D Vial: 18
Acq On : 30 May 2006 4:51 pm Operator:
Sample : 56637-5 Inst : Instrumen
Misc : 5000 Multiplr: 1.00
MS Integration Params: rteint.p
Quant Time: May 30 17:06 2006 Quant Results File: V805266B.RES

Method : C:\HPCHEM\1\METHODS\V805266B.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Tue May 30 08:22:32 2006
Response via : Initial Calibration



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May 31, 2006

CLIENT SAMPLE ID

Project Name: GP Old Town Mill

Project Number: 218736.002

Field Sample ID: SB-21 (0-4')

SAMPLE DATA

Lab Sample ID: 56637-9
Matrix: Solid
Percent Solid: 97
Dilution Factor: 116
Collection Date: 05/25/06
Lab Receipt Date: 05/27/06
Analysis Date: 05/30/06

ANALYTICAL RESULTS VOLATILE ORGANICS

COMPOUND	Quantitation Limit µg/kg	Result µg/kg	COMPOUND	Quantitation Limit µg/kg	Result µg/kg
Benzene	116	U	1,3-Dichloropropane	116	U
Bromobenzene	116	U	cis-1,3-Dichloropropene	116	U
Bromochloromethane	116	U	trans-1,3-Dichloropropene	116	U
Bromodichloromethane	87	U	2,2-Dichloropropane	116	U
Bromoform	87	U	1,1-Dichloropropene	116	U
Bromomethane	116	U	Ethylbenzene	116	U
n-butylbenzene	116	U	Hexachlorobutadiene	116	U
sec-butylbenzene	116	U	Isopropylbenzene	116	U
tert-butylbenzene	116	U	p-isopropyltoluene	116	U
Carbon Tetrachloride	116	U	Methylene Chloride	582	U
Chlorobenzene	116	U	Methyl-tert-butyl ether (MTBE)	116	U
Chloroethane	116	U	Naphthalene	116	U
Chloroform	87	U	n-Propylbenzene	116	U
Chloromethane	116	U	Styrene	116	U
2-Chlorotoluene	116	U	1,1,1,2-Tetrachloroethane	116	U
4-Chlorotoluene	116	U	1,1,2,2-Tetrachloroethane	87	U
Dibromochloromethane	87	U	Tetrachloroethene	116	**253**
1,2-Dibromo-3-chloropropane	116	U	Toluene	116	U
1,2-Dibromoethane	87	U	1,2,3-Trichlorobenzene	116	U
Dibromomethane	116	U	1,2,4-Trichlorobenzene	116	U
1,2-Dichlorobenzene	116	U	1,1,1-Trichloroethane	116	U
1,3-Dichlorobenzene	116	U	1,1,2-Trichloroethane	87	U
1,4-Dichlorobenzene	116	U	Trichloroethene	116	U
Dichlorodifluoromethane	116	U	Trichlorofluoromethane	116	U
1,1-Dichloroethane	116	U	1,2,3-Trichloropropane	116	U
1,2-Dichloroethane	87	U	1,2,4-Trimethylbenzene	116	U
1,1-Dichloroethene	87	U	1,3,5-Trimethylbenzene	116	U
cis-1,2-Dichloroethene	116	U	Vinyl Chloride	116	U
trans-1,2-Dichloroethene	116	U	o-Xylene	116	U
1,2-Dichloropropane	87	U	m,p-Xylene	116	U
Acetone	1160	**2170**	Diethyl ether	116	U
Carbon Disulfide	116	U	2-Hexanone	1160	U
Tetrahydrofuran	582	U	Methyl isobutyl ketone	1160	U
Methyl ethyl ketone	1160	U	Di-isopropyl ether (DIPE)	116	U
t-Butyl alcohol (TBA)	2330	U	Ethyl t-butyl ether (ETBE)	116	U
t-Amyl methyl ether (TAME)	116	U			

Surrogate Standard Recovery

d4-1,2-Dichloroethane	77	%	d8-Toluene	80	%	Bromofluorobenzene	84 %

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis was conducted according to: Test Methods for Evaluating Solid Waste, SW-846 Method 8260B.

COMMENTS: Results are expressed on a dry weight basis. Sample collection and analysis in accordance with SW-846 method 5035A.

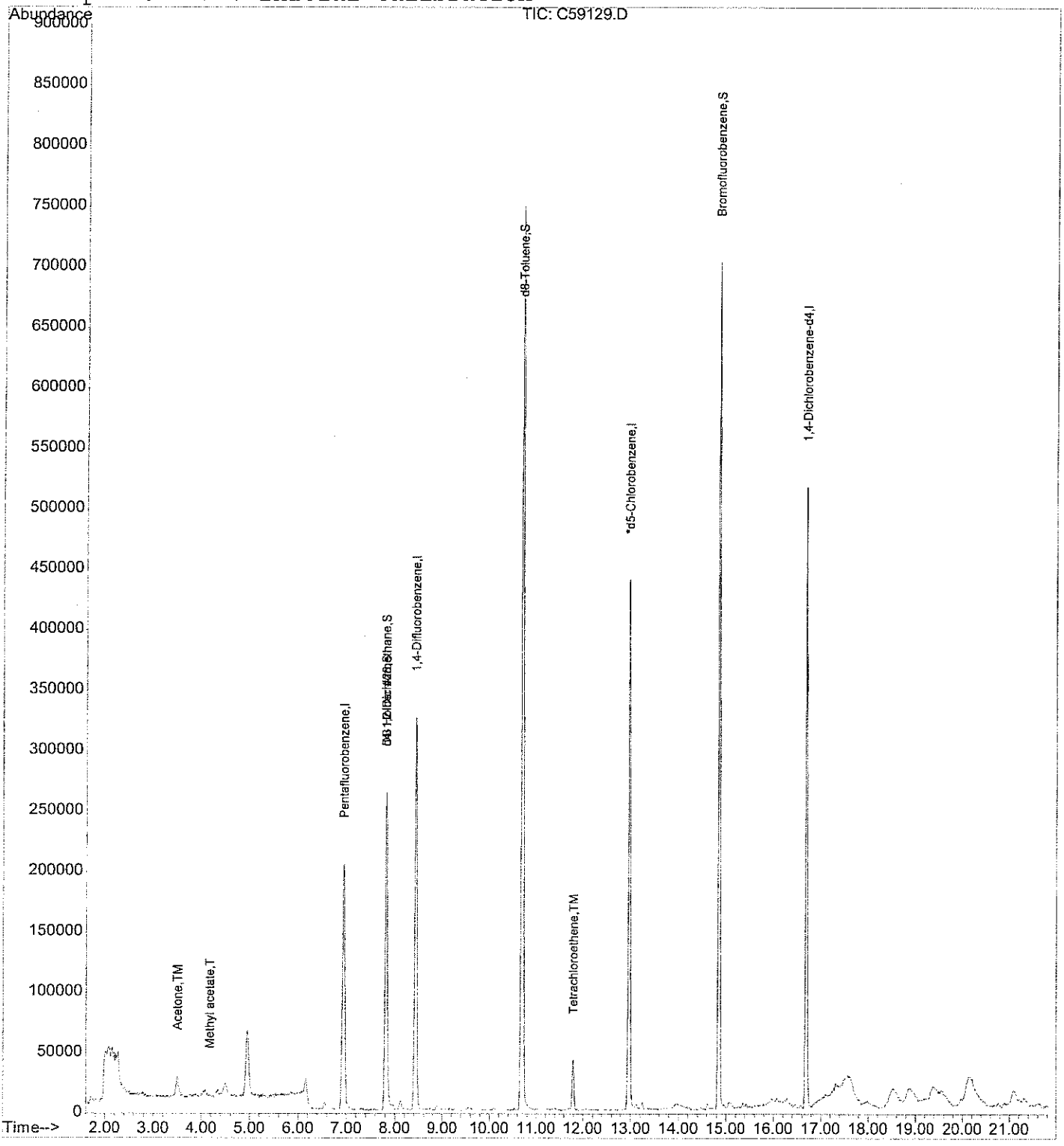
8260 full

Authorized signature

Quantitation Report

Data File : D:\HPCHEM\1\DATA\053006C\C59129.D Vial: 10
Acq On : 30 May 2006 2:46 pm Operator:
Sample : 56637-9 Inst : Instr_C
Misc : 50,5.00,SOIL Multiplr: 1.00
MS Integration Params: rteint.p
Quant Time: May 30 14:43 2006 Quant Results File: V805226C.RES

Method : D:\HPCHEM\1\METHODS\V805226C.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Mon May 22 14:18:09 2006
Response via : Initial Calibration



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May 31, 2006

CLIENT SAMPLE ID

Project Name: GP Old Town Mill

Project Number: 218736.002

Field Sample ID: SB-5 (0-4')

SAMPLE DATA

Lab Sample ID: 56637-10
Matrix: Solid
Percent Solid: 94
Dilution Factor: 108
Collection Date: 05/25/06
Lab Receipt Date: 05/27/06
Analysis Date: 05/30/06

ANALYTICAL RESULTS VOLATILE ORGANICS

COMPOUND	Quantitation Limit µg/kg	Result µg/kg	COMPOUND	Quantitation Limit µg/kg	Result µg/kg
Benzene	108	U	1,3-Dichloropropane	108	U
Bromobenzene	108	U	cis-1,3-Dichloropropene	108	U
Bromochloromethane	108	U	trans-1,3-Dichloropropene	108	U
Bromodichloromethane	81	U	2,2-Dichloropropane	108	U
Bromoform	81	U	1,1-Dichloropropene	108	U
Bromomethane	108	U	Ethylbenzene	108	U
n-butylbenzene	108	U	Hexachlorobutadiene	108	U
sec-butylbenzene	108	U	Isopropylbenzene	108	U
tert-butylbenzene	108	U	p-isopropyltoluene	108	U
Carbon Tetrachloride	108	U	Methylene Chloride	542	U
Chlorobenzene	108	U	Methyl-tert-butyl ether (MTBE)	108	U
Chloroethane	108	U	Naphthalene	108	U
Chloroform	81	U	n-Propylbenzene	108	U
Chloromethane	108	U	Styrene	108	U
2-Chlorotoluene	108	U	1,1,1,2-Tetrachloroethane	108	U
4-Chlorotoluene	108	U	1,1,2,2-Tetrachloroethane	81	U
Dibromochloromethane	81	U	Tetrachloroethene	108	U
1,2-Dibromo-3-chloropropane	108	U	Toluene	108	U
1,2-Dibromoethane	81	U	1,2,3-Trichlorobenzene	108	U
Dibromomethane	108	U	1,2,4-Trichlorobenzene	108	U
1,2-Dichlorobenzene	108	U	1,1,1-Trichloroethane	108	U
1,3-Dichlorobenzene	108	U	1,1,2-Trichloroethane	81	U
1,4-Dichlorobenzene	108	U	Trichloroethene	108	U
Dichlorodifluoromethane	108	U	Trichlorofluoromethane	108	U
1,1-Dichloroethane	108	U	1,2,3-Trichloropropane	108	U
1,2-Dichloroethane	81	U	1,2,4-Trimethylbenzene	108	U
1,1-Dichloroethene	81	U	1,3,5-Trimethylbenzene	108	U
cis-1,2-Dichloroethene	108	U	Vinyl Chloride	108	U
trans-1,2-Dichloroethene	108	U	o-Xylene	108	U
1,2-Dichloropropane	81	U	m,p-Xylene	108	U
Acetone	1080	U	Diethyl ether	108	U
Carbon Disulfide	108	U	2-Hexanone	1080	U
Tetrahydrofuran	542	U	Methyl isobutyl ketone	1080	U
Methyl ethyl ketone	1080	U	Di-isopropyl ether (DIPE)	108	U
t-Butyl alcohol (TBA)	2170	U	Ethyl t-butyl ether (ETBE)	108	U
t-Amyl methyl ether (TAME)	108	U			

Surrogate Standard Recovery

d4-1,2-Dichloroethane	76	%	d8-Toluene	80	%	Bromofluorobenzene	82 %

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis was conducted according to: Test Methods for Evaluating Solid Waste, SW-846 Method 8260B.

COMMENTS: Results are expressed on a dry weight basis. Sample collection and analysis in accordance with SW-846 method 5035A.

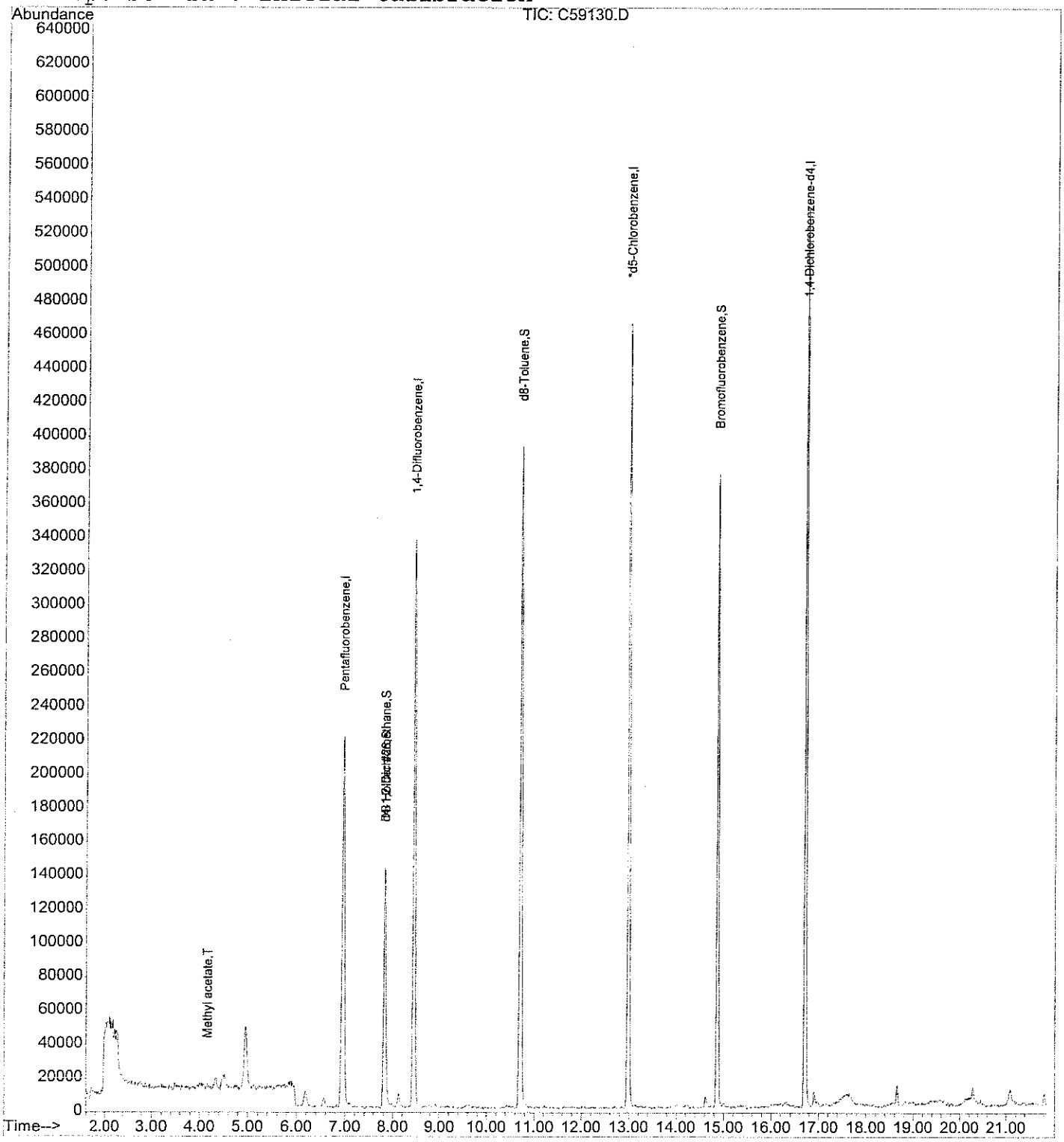
8260 full

Authorized signature

Quantitation Report

Data File : D:\HPCHEM\1\DATA\053006C\C59130.D Vial: 11
Acq On : 30 May 2006 3:21 pm Operator:
Sample : 56637-10 Inst : Instr_C
Misc : 50,5.00,SOIL Multiplr: 1.00
MS Integration Params: rteint.p
Quant Time: May 30 15:14 2006 Quant Results File: V805226C.RES

Method : D:\HPCHEM\1\METHODS\V805226C.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Mon May 22 14:18:09 2006
Response via : Initial Calibration



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Mr. Karl Kasper
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41 Hutchins Drive
Portland ME 04102

May 31, 2006

CLIENT SAMPLE ID

Project Name: GP Old Town Mill

Project Number: 218736.002

Field Sample ID: SB-6 (0-4')

SAMPLE DATA

Lab Sample ID: 56637-11
Matrix: Solid
Percent Solid: 98
Dilution Factor: 127
Collection Date: 05/25/06
Lab Receipt Date: 05/27/06
Analysis Date: 05/30/06

ANALYTICAL RESULTS VOLATILE ORGANICS

COMPOUND	Quantitation Limit µg/kg	Result µg/kg	COMPOUND	Quantitation Limit µg/kg	Result µg/kg
Benzene	127	U	1,3-Dichloropropane	127	U
Bromobenzene	127	U	cis-1,3-Dichloropropene	127	U
Bromochloromethane	127	U	trans-1,3-Dichloropropene	127	U
Bromodichloromethane	96	U	2,2-Dichloropropane	127	U
Bromoform	96	U	1,1-Dichloropropene	127	U
Bromomethane	127	U	Ethylbenzene	127	U
n-butylbenzene	127	U	Hexachlorobutadiene	127	U
sec-butylbenzene	127	U	Isopropylbenzene	127	U
tert-butylbenzene	127	U	p-isopropyltoluene	127	U
Carbon Tetrachloride	127	U	Methylene Chloride	637	U
Chlorobenzene	127	U	Methyl-tert-butyl ether (MTBE)	127	U
Chloroethane	127	U	Naphthalene	127	U
Chloroform	96	U	n-Propylbenzene	127	U
Chloromethane	127	U	Styrene	127	U
2-Chlorotoluene	127	U	1,1,1,2-Tetrachloroethane	127	U
4-Chlorotoluene	127	U	1,1,2,2-Tetrachloroethane	96	U
Dibromochloromethane	96	U	Tetrachloroethene	127	U
1,2-Dibromo-3-chloropropane	127	U	Toluene	127	U
1,2-Dibromoethane	96	U	1,2,3-Trichlorobenzene	127	U
Dibromomethane	127	U	1,2,4-Trichlorobenzene	127	U
1,2-Dichlorobenzene	127	U	1,1,1-Trichloroethane	127	U
1,3-Dichlorobenzene	127	U	1,1,2-Trichloroethane	96	U
1,4-Dichlorobenzene	127	U	Trichloroethene	127	U
Dichlorodifluoromethane	127	U	Trichlorofluoromethane	127	U
1,1-Dichloroethane	127	U	1,2,3-Trichloropropane	127	U
1,2-Dichloroethane	96	U	1,2,4-Trimethylbenzene	127	U
1,1-Dichloroethene	96	U	1,3,5-Trimethylbenzene	127	U
cis-1,2-Dichloroethene	127	U	Vinyl Chloride	127	U
trans-1,2-Dichloroethene	127	U	o-Xylene	127	U
1,2-Dichloropropane	96	U	m,p-Xylene	127	U
Acetone	1270	U	Diethyl ether	127	U
Carbon Disulfide	127	U	2-Hexanone	1270	U
Tetrahydrofuran	637	U	Methyl isobutyl ketone	1270	U
Methyl ethyl ketone	1270	U	Di-isopropyl ether (DIPE)	127	U
t-Butyl alcohol (TBA)	2550	U	Ethyl t-butyl ether (ETBE)	127	U
t-Amyl methyl ether (TAME)	127	U			

Surrogate Standard Recovery

d4-1,2-Dichloroethane	71	%	d8-Toluene	78	%	Bromofluorobenzene	79 %

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis was conducted according to: Test Methods for Evaluating Solid Waste, SW-846 Method 8260B.

COMMENTS: Results are expressed on a dry weight basis. Sample collection and analysis in accordance with SW-846 method 5035A.

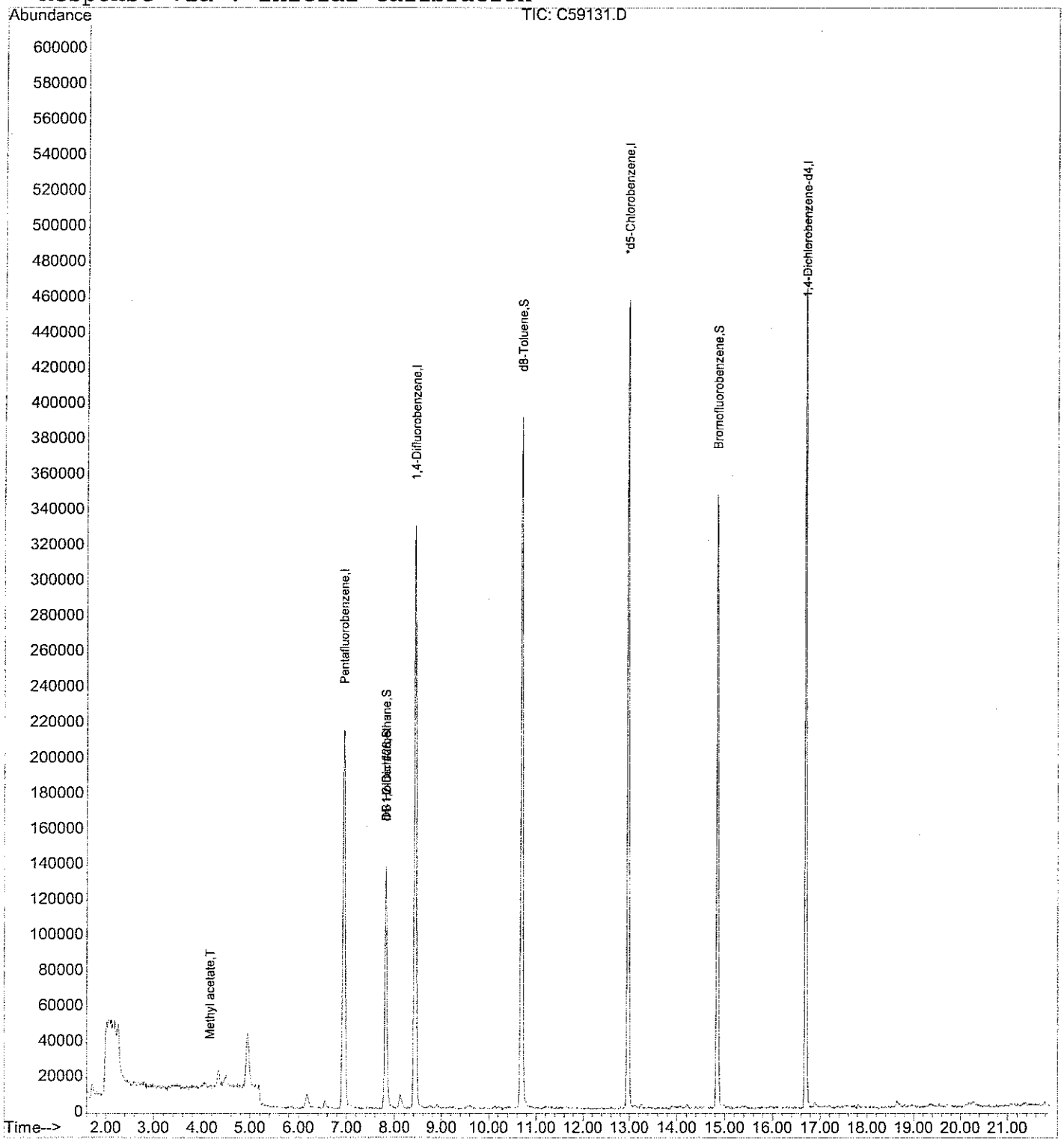
8260 full

Authorized signature

Quantitation Report

Data File : D:\HPCHEM\1\DATA\053006C\C59131.D Vial: 12
Acq On : 30 May 2006 3:55 pm Operator:
Sample : 56637-11 Inst : Instr_C
Misc : 50,5.00,SOIL Multiplr: 1.00
MS Integration Params: rteint.p
Quant Time: May 30 15:44 2006 Quant Results File: V805226C.RES

Method : D:\HPCHEM\1\METHODS\V805226C.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Mon May 22 14:18:09 2006
Response via : Initial Calibration



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Portland ME 04102

May 31, 2006

CLIENT SAMPLE ID

Project Name: GP Old Town Mill

Project Number: 218736.002

Field Sample ID: SB-31 (0-4)

SAMPLE DATA

Lab Sample ID: 56637-14
Matrix: Solid
Percent Solid: 63
Dilution Factor: 193
Collection Date: 05/25/06
Lab Receipt Date: 05/27/06
Analysis Date: 05/30/06

ANALYTICAL RESULTS VOLATILE ORGANICS

COMPOUND	Quantitation Limit µg/kg	Result µg/kg	COMPOUND	Quantitation Limit µg/kg	Result µg/kg
Benzene	193	U	1,3-Dichloropropane	193	U
Bromobenzene	193	U	cis-1,3-Dichloropropene	193	U
Bromochloromethane	193	U	trans-1,3-Dichloropropene	193	U
Bromodichloromethane	144	U	2,2-Dichloropropane	193	U
Bromoform	144	U	1,1-Dichloropropene	193	U
Bromomethane	193	U	Ethylbenzene	193	U
n-butylbenzene	193	U	Hexachlorobutadiene	193	U
sec-butylbenzene	193	U	Isopropylbenzene	193	U
tert-butylbenzene	193	U	p-isopropyltoluene	193	U
Carbon Tetrachloride	193	U	Methylene Chloride	963	U
Chlorobenzene	193	U	Methyl-tert-butyl ether (MTBE)	193	U
Chloroethane	193	U	Naphthalene	193	U
Chloroform	144	U	n-Propylbenzene	193	U
Chloromethane	193	U	Styrene	193	U
2-Chlorotoluene	193	U	1,1,1,2-Tetrachloroethane	193	U
4-Chlorotoluene	193	U	1,1,2,2-Tetrachloroethane	144	U
Dibromochloromethane	144	U	Tetrachloroethene	193	U
1,2-Dibromo-3-chloropropane	193	U	Toluene	193	U
1,2-Dibromoethane	144	U	1,2,3-Trichlorobenzene	193	U
Dibromomethane	193	U	1,2,4-Trichlorobenzene	193	U
1,2-Dichlorobenzene	193	U	1,1,1-Trichloroethane	193	U
1,3-Dichlorobenzene	193	U	1,1,2-Trichloroethane	144	U
1,4-Dichlorobenzene	193	U	Trichloroethene	193	U
Dichlorodifluoromethane	193	U	Trichlorofluoromethane	193	U
1,1-Dichloroethane	193	U	1,2,3-Trichloropropane	193	U
1,2-Dichloroethane	144	U	1,2,4-Trimethylbenzene	193	U
1,1-Dichloroethene	144	U	1,3,5-Trimethylbenzene	193	U
cis-1,2-Dichloroethene	193	U	Vinyl Chloride	193	U
trans-1,2-Dichloroethene	193	U	o-Xylene	193	U
1,2-Dichloropropane	144	U	m,p-Xylene	193	U
Acetone	1930	U	Diethyl ether	193	U
Carbon Disulfide	193	U	2-Hexanone	1930	U
Tetrahydrofuran	963	U	Methyl isobutyl ketone	1930	U
Methyl ethyl ketone	1930	U	Di-isopropyl ether (DIPE)	193	U
t-Butyl alcohol (TBA)	3850	U	Ethyl t-butyl ether (ETBE)	193	U
t-Amyl methyl ether (TAME)	193	U			

Surrogate Standard Recovery

d4-1,2-Dichloroethane	64 * %	d8-Toluene	69 * %	Bromofluorobenzene	67 * %

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis was conducted according to: Test Methods for Evaluating Solid Waste, SW-846 Method 8260B.

COMMENTS: Results are expressed on a dry weight basis. Sample collection and analysis in accordance with SW-846 method 5035A. * Surrogate recovery outside laboratory acceptance criteria. Sample was reanalyzed with similar results.

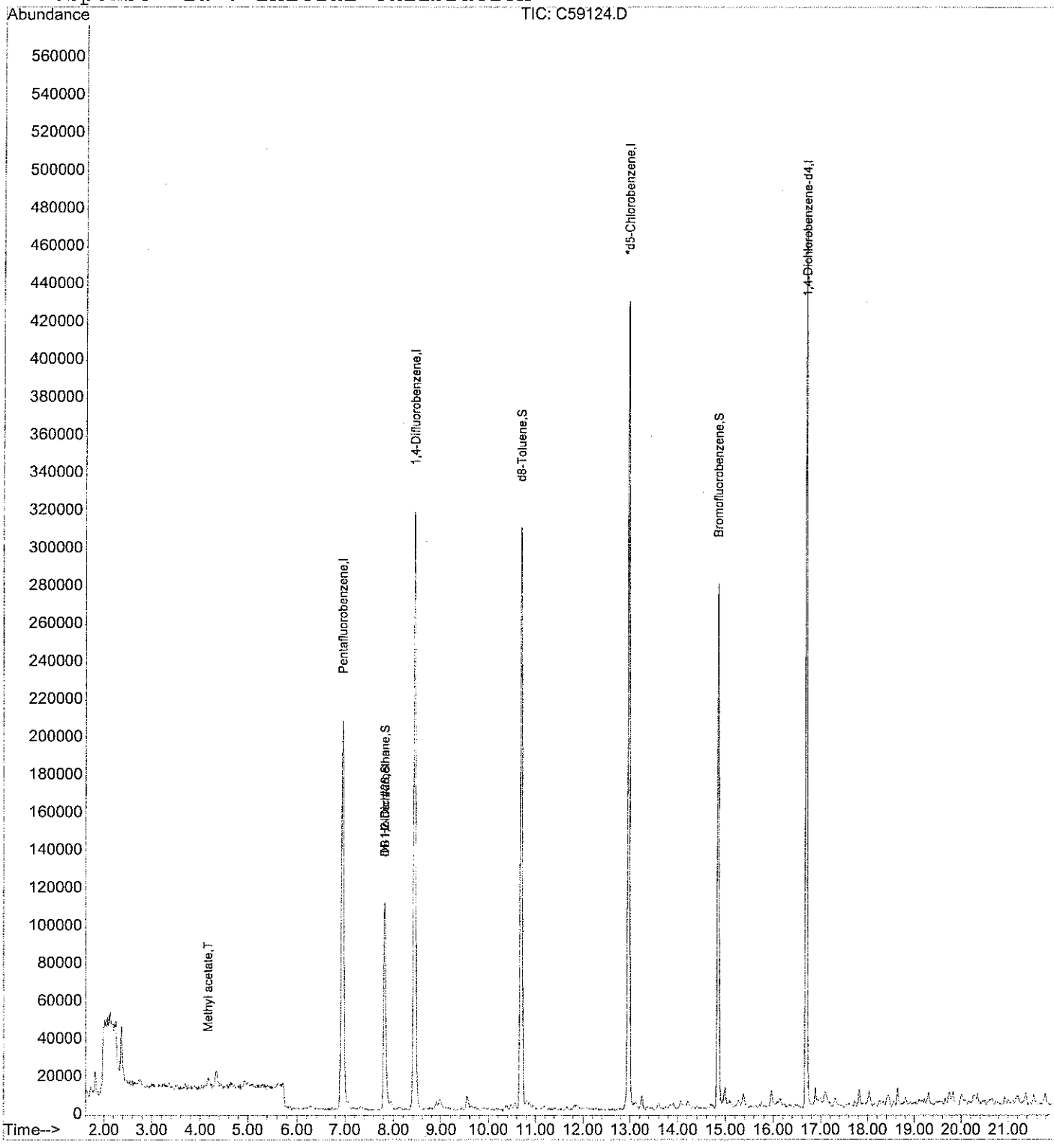
8260 full

Authorized signature

Quantitation Report

Data File : D:\HPCHEM\1\DATA\053006C\C59124.D Vial: 16
Acq On : 30 May 2006 12:03 pm Operator:
Sample : 56637-14 Inst : Instr_C
Misc : 50,4.10,SOIL Multiplr: 1.00
MS Integration Params: rteint.p
Quant Time: May 30 12:07 2006 Quant Results File: V805226C.RES

Method : D:\HPCHEM\1\METHODS\V805226C.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Mon May 22 14:18:09 2006
Response via : Initial Calibration



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May 31, 2006

CLIENT SAMPLE ID

Project Name: GP Old Town Mill

Project Number: 218736.002

Field Sample ID: SB-30 (0-4)

SAMPLE DATA

Lab Sample ID: 56637-15
Matrix: Solid
Percent Solid: 81
Dilution Factor: 200
Collection Date: 05/25/06
Lab Receipt Date: 05/27/06
Analysis Date: 05/30/06

ANALYTICAL RESULTS VOLATILE ORGANICS

COMPOUND	Quantitation Limit µg/kg	Result µg/kg	COMPOUND	Quantitation Limit µg/kg	Result µg/kg
Benzene	200	U	1,3-Dichloropropane	200	U
Bromobenzene	200	U	cis-1,3-Dichloropropene	200	U
Bromochloromethane	200	U	trans-1,3-Dichloropropene	200	U
Bromodichloromethane	150	U	2,2-Dichloropropane	200	U
Bromoform	150	U	1,1-Dichloropropene	200	U
Bromomethane	200	U	Ethylbenzene	200	U
n-butylbenzene	200	U	Hexachlorobutadiene	200	U
sec-butylbenzene	200	U	Isopropylbenzene	200	U
tert-butylbenzene	200	U	p-isopropyltoluene	200	U
Carbon Tetrachloride	200	U	Methylene Chloride	998	U
Chlorobenzene	200	U	Methyl-tert-butyl ether (MTBE)	200	U
Chloroethane	200	U	Naphthalene	200	U
Chloroform	150	U	n-Propylbenzene	200	U
Chloromethane	200	U	Styrene	200	U
2-Chlorotoluene	200	U	1,1,1,2-Tetrachloroethane	200	U
4-Chlorotoluene	200	U	1,1,2,2-Tetrachloroethane	150	U
Dibromochloromethane	150	U	Tetrachloroethene	200	U
1,2-Dibromo-3-chloropropane	200	U	Toluene	200	U
1,2-Dibromoethane	150	U	1,2,3-Trichlorobenzene	200	U
Dibromomethane	200	U	1,2,4-Trichlorobenzene	200	U
1,2-Dichlorobenzene	200	U	1,1,1-Trichloroethane	200	U
1,3-Dichlorobenzene	200	U	1,1,2-Trichloroethane	150	U
1,4-Dichlorobenzene	200	U	Trichloroethene	200	U
Dichlorodifluoromethane	200	U	Trichlorofluoromethane	200	U
1,1-Dichloroethane	200	U	1,2,3-Trichloropropane	200	U
1,2-Dichloroethane	150	U	1,2,4-Trimethylbenzene	200	U
1,1-Dichloroethene	150	U	1,3,5-Trimethylbenzene	200	U
cis-1,2-Dichloroethene	200	U	Vinyl Chloride	200	U
trans-1,2-Dichloroethene	200	U	o-Xylene	200	U
1,2-Dichloropropane	150	U	m,p-Xylene	200	U
Acetone	2000	U	Diethyl ether	200	U
Carbon Disulfide	200	U	2-Hexanone	2000	U
Tetrahydrofuran	998	U	Methyl isobutyl ketone	2000	U
Methyl ethyl ketone	2000	U	Di-isopropyl ether (DIPE)	200	U
t-Butyl alcohol (TBA)	3990	U	Ethyl t-butyl ether (ETBE)	200	U
t-Amyl methyl ether (TAME)	200	U			

Surrogate Standard Recovery

d4-1,2-Dichloroethane	99 %	d8-Toluene	103 %	Bromofluorobenzene	103 %

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis was conducted according to: Test Methods for Evaluating Solid Waste, SW-846 Method 8260B.

COMMENTS: Results are expressed on a dry weight basis. Sample collection and analysis in accordance with SW-846 method 5035. Sample did not meet method acceptance criteria for 1:1 soil to methanol ratio.

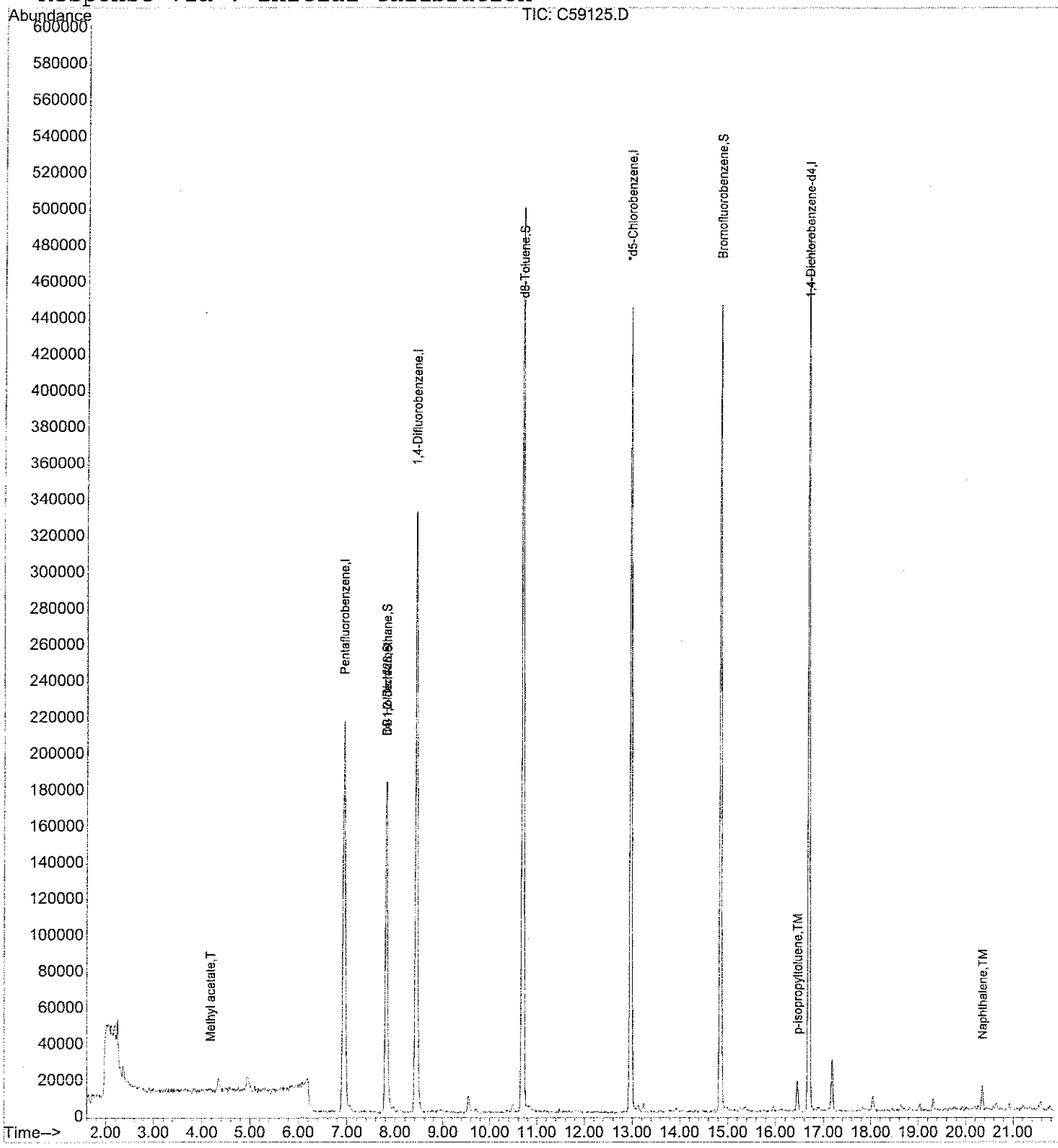
8260 full

Authorized signature

Quantitation Report

Data File : D:\HPCHEM\1\DATA\053006C\C59125.D Vial: 6
Acq On : 30 May 2006 12:36 pm Operator:
Sample : 56637-15 Inst : Instr_C
Misc : 50,3.08,SOIL Multiplr: 1.00
MS Integration Params: rteint.p
Quant Time: May 30 12:38 2006 Quant Results File: V805226C.RES

Method : D:\HPCHEM\1\METHODS\V805226C.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Mon May 22 14:18:09 2006
Response via : Initial Calibration



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May 31, 2006

CLIENT SAMPLE ID

Project Name: GP Old Town Mill

Project Number: 218736.002

Field Sample ID: SB-38 (0-4)

SAMPLE DATA

Lab Sample ID: 56637-16
Matrix: Solid
Percent Solid: 94
Dilution Factor: 107
Collection Date: 05/25/06
Lab Receipt Date: 05/27/06
Analysis Date: 05/30/06

ANALYTICAL RESULTS VOLATILE ORGANICS

COMPOUND	Quantitation Limit µg/kg	Result µg/kg	COMPOUND	Quantitation Limit µg/kg	Result µg/kg
Benzene	107	U	1,3-Dichloropropane	107	U
Bromobenzene	107	U	cis-1,3-Dichloropropene	107	U
Bromochloromethane	107	U	trans-1,3-Dichloropropene	107	U
Bromodichloromethane	80	U	2,2-Dichloropropane	107	U
Bromoform	80	U	1,1-Dichloropropene	107	U
Bromomethane	107	U	Ethylbenzene	107	U
n-butylbenzene	107	U	Hexachlorobutadiene	107	U
sec-butylbenzene	107	U	Isopropylbenzene	107	U
tert-butylbenzene	107	U	p-isopropyltoluene	107	**272**
Carbon Tetrachloride	107	U	Methylene Chloride	537	U
Chlorobenzene	107	U	Methyl-tert-butyl ether (MTBE)	107	U
Chloroethane	107	U	Naphthalene	107	**9400**
Chloroform	80	U	n-Propylbenzene	107	U
Chloromethane	107	U	Styrene	107	**292**
2-Chlorotoluene	107	U	1,1,1,2-Tetrachloroethane	107	U
4-Chlorotoluene	107	U	1,1,2,2-Tetrachloroethane	80	U
Dibromochloromethane	80	U	Tetrachloroethene	107	U
1,2-Dibromo-3-chloropropane	107	U	Toluene	107	**120**
1,2-Dibromoethane	80	U	1,2,3-Trichlorobenzene	107	U
Dibromomethane	107	U	1,2,4-Trichlorobenzene	107	U
1,2-Dichlorobenzene	107	U	1,1,1-Trichloroethane	107	U
1,3-Dichlorobenzene	107	U	1,1,2-Trichloroethane	80	U
1,4-Dichlorobenzene	107	U	Trichloroethene	107	U
Dichlorodifluoromethane	107	U	Trichlorofluoromethane	107	U
1,1-Dichloroethane	107	U	1,2,3-Trichloropropane	107	U
1,2-Dichloroethane	80	U	1,2,4-Trimethylbenzene	107	**1510**
1,1-Dichloroethene	80	U	1,3,5-Trimethylbenzene	107	**1150**
cis-1,2-Dichloroethene	107	U	Vinyl Chloride	107	U
trans-1,2-Dichloroethene	107	U	o-Xylene	107	**333**
1,2-Dichloropropane	80	U	m,p-Xylene	107	**347**
Acetone	1070	U	Diethyl ether	107	U
Carbon Disulfide	107	U	2-Hexanone	1070	U
Tetrahydrofuran	537	U	Methyl isobutyl ketone	1070	U
Methyl ethyl ketone	1070	U	Di-isopropyl ether (DIPE)	107	U
t-Butyl alcohol (TBA)	2150	U	Ethyl t-butyl ether (ETBE)	107	U
t-Amyl methyl ether (TAME)	107	U			

Surrogate Standard Recovery

d4-1,2-Dichloroethane	78	%	d8-Toluene	79	%	Bromofluorobenzene	84 %

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis was conducted according to: Test Methods for Evaluating Solid Waste, SW-846 Method 8260B.

COMMENTS: Results are expressed on a dry weight basis. Sample collection and analysis in accordance with SW-846 method 5035A.

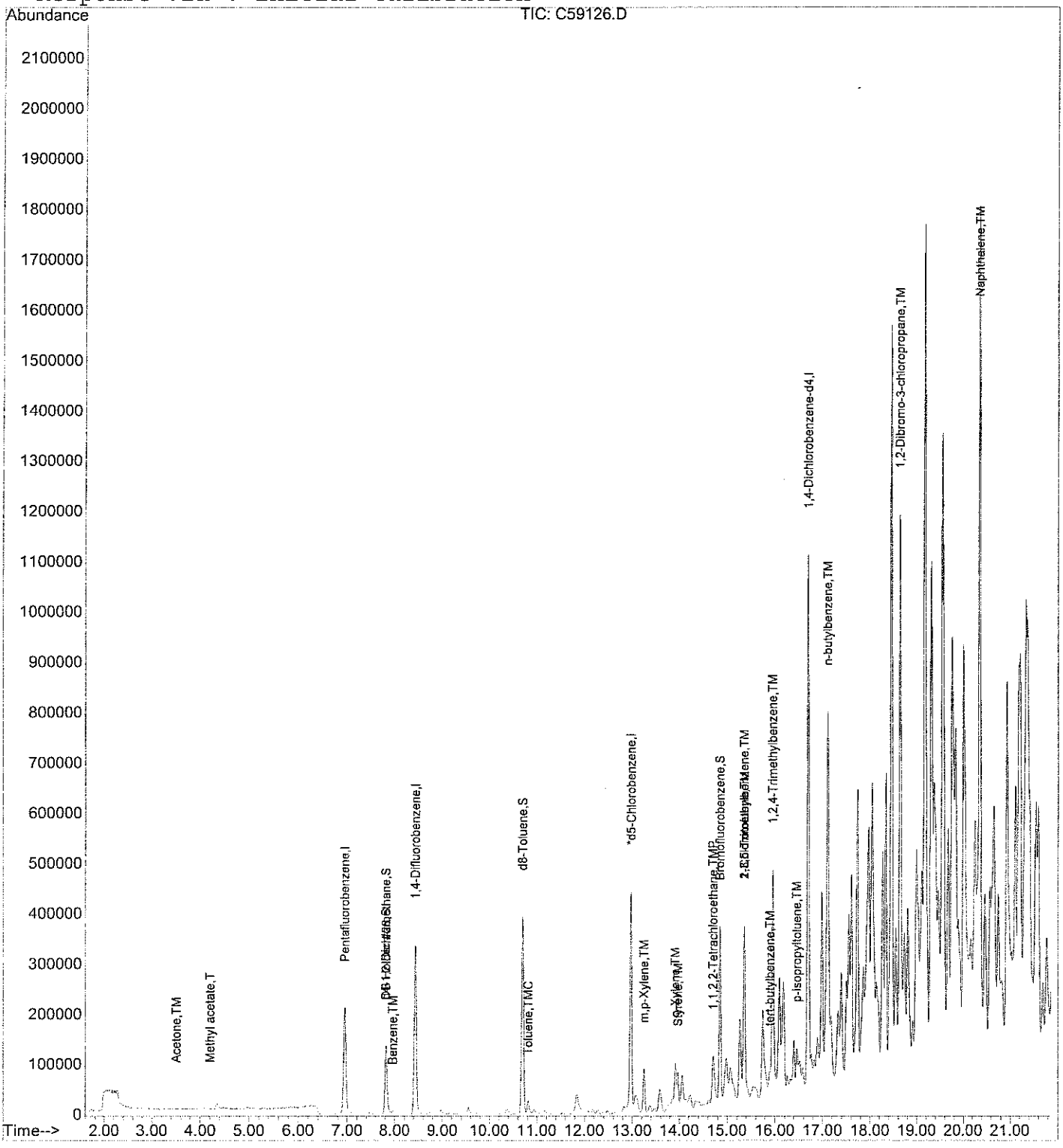
8260 full

Authorized signature

Quantitation Report

Data File : D:\HPCHEM\1\DATA\053006C\C59126.D Vial: 7
Acq On : 30 May 2006 1:14 pm Operator:
Sample : 56637-16 Inst : Instr_C
Misc : 50,4.98,SOIL Multiplr: 1.00
MS Integration Params: rteint.p
Quant Time: May 30 13:09 2006 Quant Results File: V805226C.RES

Method : D:\HPCHEM\1\METHODS\V805226C.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Mon May 22 14:18:09 2006
Response via : Initial Calibration



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May 31, 2006

CLIENT SAMPLE ID

Project Name: GP Old Town Mill

Project Number: 218736.002

Field Sample ID: SB-37 (0-4)

SAMPLE DATA

Lab Sample ID: 56637-17 RR
Matrix: Solid
Percent Solid: 93
Dilution Factor: 113
Collection Date: 05/25/06
Lab Receipt Date: 05/27/06
Analysis Date: 05/30/06

ANALYTICAL RESULTS VOLATILE ORGANICS

COMPOUND	Quantitation Limit µg/kg	Result µg/kg	COMPOUND	Quantitation Limit µg/kg	Result µg/kg
Benzene	113	U	1,3-Dichloropropane	113	U
Bromobenzene	113	U	cis-1,3-Dichloropropene	113	U
Bromochloromethane	113	U	trans-1,3-Dichloropropene	113	U
Bromodichloromethane	84	U	2,2-Dichloropropane	113	U
Bromoform	84	U	1,1-Dichloropropene	113	U
Bromomethane	113	U	Ethylbenzene	113	U
n-butylbenzene	113	U	Hexachlorobutadiene	113	U
sec-butylbenzene	113	U	Isopropylbenzene	113	U
tert-butylbenzene	113	U	p-isopropyltoluene	113	U
Carbon Tetrachloride	113	U	Methylene Chloride	563	U
Chlorobenzene	113	U	Methyl-tert-butyl ether (MTBE)	113	U
Chloroethane	113	U	Naphthalene	113	U
Chloroform	84	U	n-Propylbenzene	113	U
Chloromethane	113	U	Styrene	113	U
2-Chlorotoluene	113	U	1,1,1,2-Tetrachloroethane	113	U
4-Chlorotoluene	113	U	1,1,2,2-Tetrachloroethane	84	U
Dibromochloromethane	84	U	Tetrachloroethene	113	U
1,2-Dibromo-3-chloropropane	113	U	Toluene	113	U
1,2-Dibromoethane	84	U	1,2,3-Trichlorobenzene	113	U
Dibromomethane	113	U	1,2,4-Trichlorobenzene	113	U
1,2-Dichlorobenzene	113	U	1,1,1-Trichloroethane	113	U
1,3-Dichlorobenzene	113	U	1,1,2-Trichloroethane	84	U
1,4-Dichlorobenzene	113	U	Trichloroethene	113	U
Dichlorodifluoromethane	113	U	Trichlorofluoromethane	113	U
1,1-Dichloroethane	113	U	1,2,3-Trichloropropane	113	U
1,2-Dichloroethane	84	U	1,2,4-Trimethylbenzene	113	U
1,1-Dichloroethene	84	U	1,3,5-Trimethylbenzene	113	U
cis-1,2-Dichloroethene	113	U	Vinyl Chloride	113	U
trans-1,2-Dichloroethene	113	U	o-Xylene	113	U
1,2-Dichloropropane	84	U	m,p-Xylene	113	U
Acetone	1130	U	Diethyl ether	113	U
Carbon Disulfide	113	U	2-Hexanone	1130	U
Tetrahydrofuran	563	U	Methyl isobutyl ketone	1130	U
Methyl ethyl ketone	1130	U	Di-isopropyl ether (DIPE)	113	U
t-Butyl alcohol (TBA)	2250	U	Ethyl t-butyl ether (ETBE)	113	U
t-Amyl methyl ether (TAME)	113	U			

Surrogate Standard Recovery

d4-1,2-Dichloroethane	73 %	d8-Toluene	78 %	Bromofluorobenzene	82 %

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis was conducted according to: Test Methods for Evaluating Solid Waste, SW-846 Method 8260B.

COMMENTS: Results are expressed on a dry weight basis. Sample collection and analysis in accordance with SW-846 method 5035A.

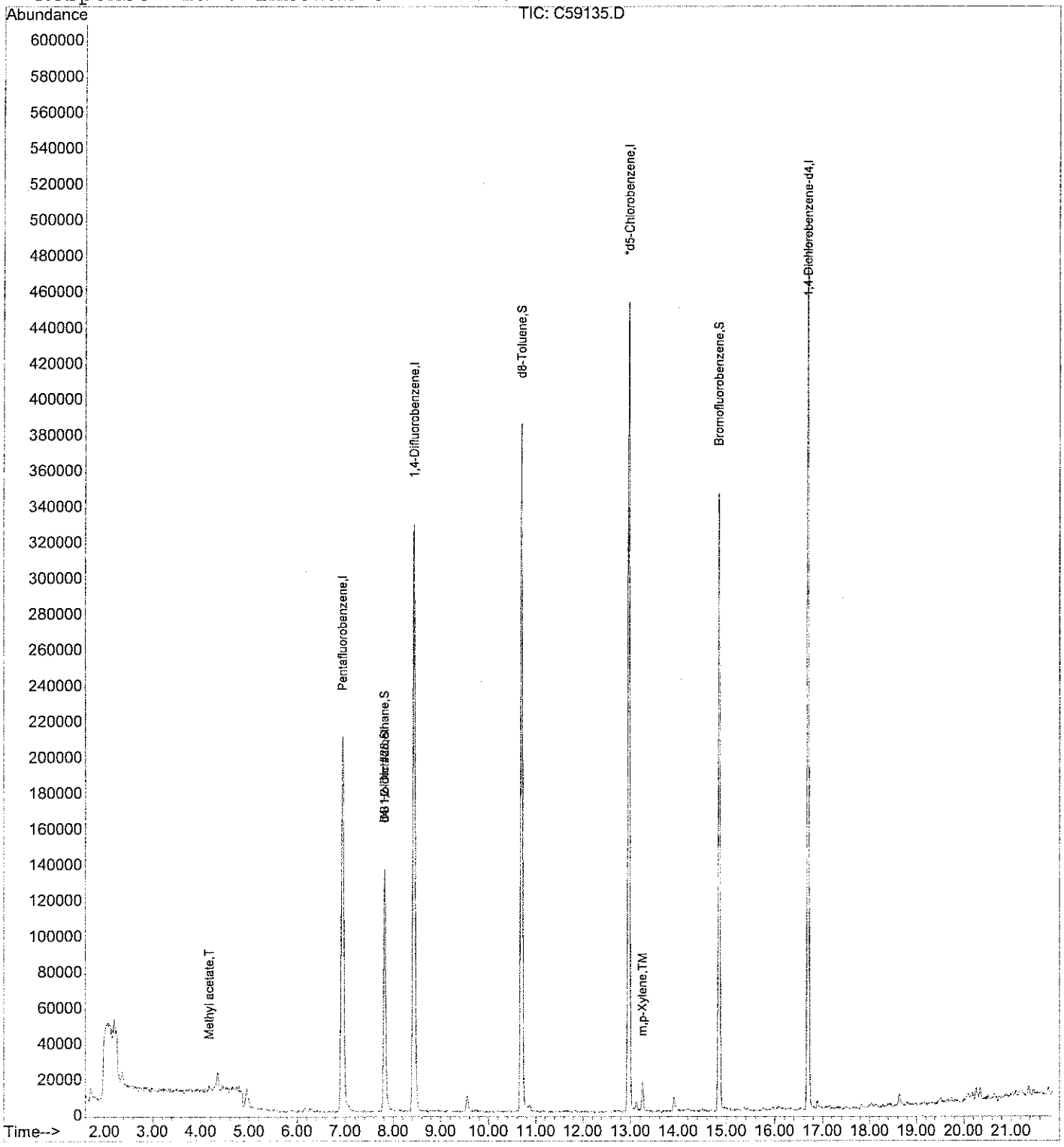
8260 full

Authorized signature

Quantitation Report

Data File : D:\HPCHEM\1\DATA\053006C\C59135.D Vial: 16
Acq On : 30 May 2006 6:33 pm Operator:
Sample : 56637-17, RR Inst : Instr_C
Misc : 50,4.78,SOIL Multiplr: 1.00
MS Integration Params: rteint.p
Quant Time: May 30 18:29 2006 Quant Results File: V805226C.RES

Method : D:\HPCHEM\1\METHODS\V805226C.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Mon May 22 14:18:09 2006
Response via : Initial Calibration



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Mr. Karl Kasper
Woodard & Curran
41 Hutchins Drive
Portland ME 04102

May 31, 2006

CLIENT SAMPLE ID

Project Name: GP Old Town Mill

Project Number: 218736.002

Field Sample ID: SB-15 (0-4)

SAMPLE DATA

Lab Sample ID: 56637-18 RR
Matrix: Solid
Percent Solid: 88
Dilution Factor: 176
Collection Date: 05/25/06
Lab Receipt Date: 05/27/06
Analysis Date: 05/30/06

ANALYTICAL RESULTS VOLATILE ORGANICS

COMPOUND	Quantitation Limit µg/kg	Result µg/kg	COMPOUND	Quantitation Limit µg/kg	Result µg/kg
Benzene	176	U	1,3-Dichloropropane	176	U
Bromobenzene	176	U	cis-1,3-Dichloropropene	176	U
Bromochloromethane	176	U	trans-1,3-Dichloropropene	176	U
Bromodichloromethane	132	U	2,2-Dichloropropane	176	U
Bromoform	132	U	1,1-Dichloropropene	176	U
Bromomethane	176	U	Ethylbenzene	176	U
n-butylbenzene	176	U	Hexachlorobutadiene	176	U
sec-butylbenzene	176	U	Isopropylbenzene	176	U
tert-butylbenzene	176	U	p-isopropyltoluene	176	U
Carbon Tetrachloride	176	U	Methylene Chloride	882	U
Chlorobenzene	176	U	Methyl-tert-butyl ether (MTBE)	176	U
Chloroethane	176	U	Naphthalene	176	U
Chloroform	132	U	n-Propylbenzene	176	U
Chloromethane	176	U	Styrene	176	U
2-Chlorotoluene	176	U	1,1,1,2-Tetrachloroethane	176	U
4-Chlorotoluene	176	U	1,1,2,2-Tetrachloroethane	132	U
Dibromochloromethane	132	U	Tetrachloroethene	176	U
1,2-Dibromo-3-chloropropane	176	U	Toluene	176	U
1,2-Dibromoethane	132	U	1,2,3-Trichlorobenzene	176	U
Dibromomethane	176	U	1,2,4-Trichlorobenzene	176	U
1,2-Dichlorobenzene	176	U	1,1,1-Trichloroethane	176	U
1,3-Dichlorobenzene	176	U	1,1,2-Trichloroethane	132	U
1,4-Dichlorobenzene	176	U	Trichloroethene	176	U
Dichlorodifluoromethane	176	U	Trichlorofluoromethane	176	U
1,1-Dichloroethane	176	U	1,2,3-Trichloropropane	176	U
1,2-Dichloroethane	132	U	1,2,4-Trimethylbenzene	176	U
1,1-Dichloroethene	132	U	1,3,5-Trimethylbenzene	176	U
cis-1,2-Dichloroethene	176	U	Vinyl Chloride	176	U
trans-1,2-Dichloroethene	176	U	o-Xylene	176	U
1,2-Dichloropropane	132	U	m,p-Xylene	176	U
Acetone	1760	U	Diethyl ether	176	U
Carbon Disulfide	176	U	2-Hexanone	1760	U
Tetrahydrofuran	882	U	Methyl isobutyl ketone	1760	U
Methyl ethyl ketone	1760	U	Di-isopropyl ether (DIPE)	176	U
t-Butyl alcohol (TBA)	3530	U	Ethyl t-butyl ether (ETBE)	176	U
t-Amyl methyl ether (TAME)	176	U			

Surrogate Standard Recovery

d4-1,2-Dichloroethane	71 %	d8-Toluene	70 %	Bromofluorobenzene	72 %

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis was conducted according to: Test Methods for Evaluating Solid Waste, SW-846 Method 8260B.

COMMENTS: Results are expressed on a dry weight basis. Sample collection and analysis in accordance with SW-846 method 5035A.

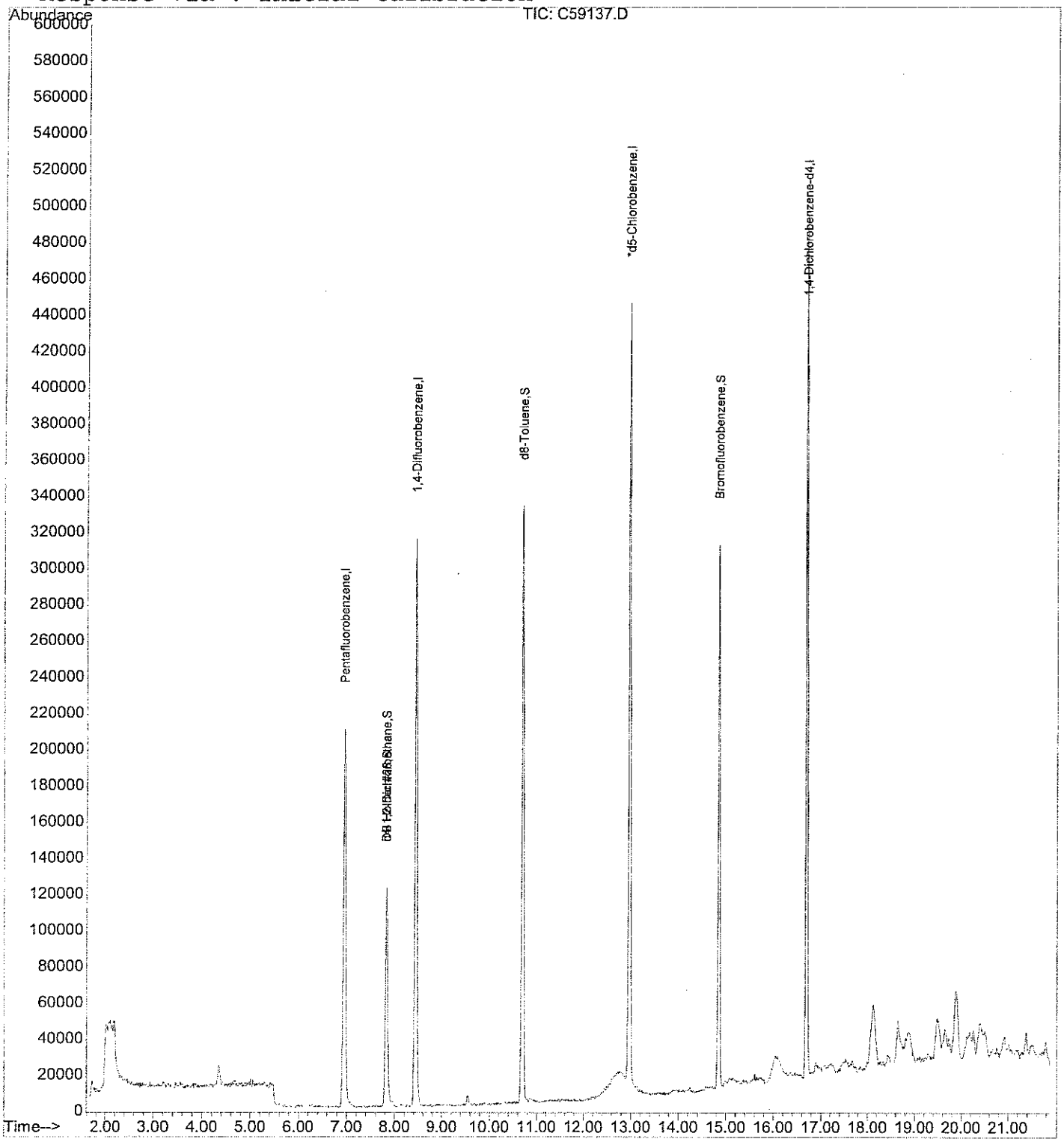
8260 full

Authorized signature Milena Tull

Quantitation Report

Data File : D:\HPCHEM\1\DATA\053006C\C59137.D Vial: 2
Acq On : 30 May 2006 7:10 pm Operator:
Sample : 56637-18, RR,, 10ML FV Inst : Instr_C
Misc : 50,5.00,SOIL,,10ML FV Multiplr: 1.00
MS Integration Params: rteint.p
Quant Time: May 30 19:30 2006 Quant Results File: V805226C.RES

Method : D:\HPCHEM\1\METHODS\V805226C.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Mon May 22 14:18:09 2006
Response via : Initial Calibration



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May 31, 2006

CLIENT SAMPLE ID

Project Name: GP Old Town Mill

Project Number: 218736.002

Field Sample ID: SB-41 (0-4)

SAMPLE DATA

Lab Sample ID: 56637-19
Matrix: Solid
Percent Solid: 91
Dilution Factor: 126
Collection Date: 05/26/06
Lab Receipt Date: 05/27/06
Analysis Date: 05/30/06

ANALYTICAL RESULTS VOLATILE ORGANICS

COMPOUND	Quantitation Limit µg/kg	Result µg/kg	COMPOUND	Quantitation Limit µg/kg	Result µg/kg
Benzene	126	U	1,3-Dichloropropane	126	U
Bromobenzene	126	U	cis-1,3-Dichloropropene	126	U
Bromochloromethane	126	U	trans-1,3-Dichloropropene	126	U
Bromodichloromethane	95	U	2,2-Dichloropropane	126	U
Bromoform	95	U	1,1-Dichloropropene	126	U
Bromomethane	126	U	Ethylbenzene	126	U
n-butylbenzene	126	U	Hexachlorobutadiene	126	U
sec-butylbenzene	126	U	Isopropylbenzene	126	U
tert-butylbenzene	126	U	p-isopropyltoluene	126	U
Carbon Tetrachloride	126	U	Methylene Chloride	632	U
Chlorobenzene	126	U	Methyl-tert-butyl ether (MTBE)	126	U
Chloroethane	126	U	Naphthalene	126	U
Chloroform	95	U	n-Propylbenzene	126	U
Chloromethane	126	U	Styrene	126	U
2-Chlorotoluene	126	U	1,1,1,2-Tetrachloroethane	126	U
4-Chlorotoluene	126	U	1,1,2,2-Tetrachloroethane	95	U
Dibromochloromethane	95	U	Tetrachloroethene	126	U
1,2-Dibromo-3-chloropropane	126	U	Toluene	126	U
1,2-Dibromoethane	95	U	1,2,3-Trichlorobenzene	126	U
Dibromomethane	126	U	1,2,4-Trichlorobenzene	126	U
1,2-Dichlorobenzene	126	U	1,1,1-Trichloroethane	126	U
1,3-Dichlorobenzene	126	U	1,1,2-Trichloroethane	95	U
1,4-Dichlorobenzene	126	U	Trichloroethene	126	U
Dichlorodifluoromethane	126	U	Trichlorofluoromethane	126	U
1,1-Dichloroethane	126	U	1,2,3-Trichloropropane	126	U
1,2-Dichloroethane	95	U	1,2,4-Trimethylbenzene	126	U
1,1-Dichloroethene	95	U	1,3,5-Trimethylbenzene	126	U
cis-1,2-Dichloroethene	126	U	Vinyl Chloride	126	U
trans-1,2-Dichloroethene	126	U	o-Xylene	126	U
1,2-Dichloropropane	95	U	m,p-Xylene	126	U
Acetone	1260	U	Diethyl ether	126	U
Carbon Disulfide	126	U	2-Hexanone	1260	U
Tetrahydrofuran	632	U	Methyl isobutyl ketone	1260	U
Methyl ethyl ketone	1260	U	Di-isopropyl ether (DIPE)	126	U
t-Butyl alcohol (TBA)	2530	U	Ethyl t-butyl ether (ETBE)	126	U
t-Amyl methyl ether (TAME)	126	U			

Surrogate Standard Recovery

d4-1,2-Dichloroethane	98 %	d8-Toluene	95 %	Bromofluorobenzene	93 %

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis was conducted according to: Test Methods for Evaluating Solid Waste, SW-846 Method 8260B.

COMMENTS: Results are expressed on a dry weight basis. Sample collection and analysis in accordance with SW-846 method 5035A.

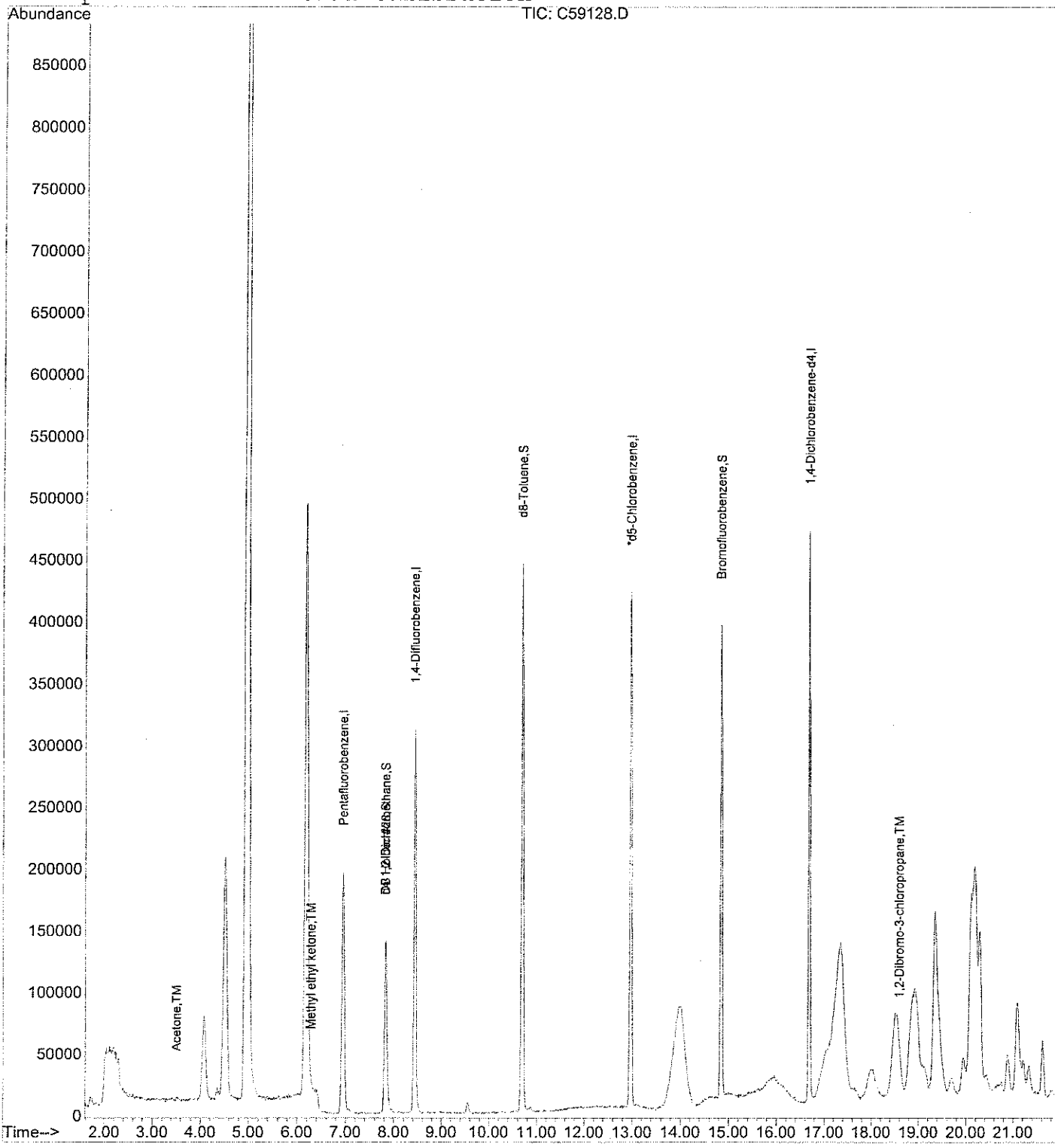
8260 full

Authorized signature

Quantitation Report

Data File : D:\HPCHEM\1\DATA\053006C\C59128.D Vial: 9
Acq On : 30 May 2006 2:17 pm Operator:
Sample : 56637-19 Inst : Instr_C
Misc : 50,4.37,SOIL Multiplr: 1.00
MS Integration Params: rteint.p
Quant Time: May 30 14:12 2006 Quant Results File: V805226C.RES

Method : D:\HPCHEM\1\METHODS\V805226C.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Mon May 22 14:18:09 2006
Response via : Initial Calibration



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May 31, 2006

CLIENT SAMPLE ID

Project Name: GP Old Town Mill

Project Number: 218736.002

Field Sample ID: SB-15

SAMPLE DATA

Lab Sample ID: 56637-20
Matrix: Aqueous
Percent Solid: N/A
Dilution Factor: 1
Collection Date: 05/26/06
Lab Receipt Date: 05/27/06
Analysis Date: 05/30/06

ANALYTICAL RESULTS VOLATILE ORGANICS

COMPOUND	Quantitation Limit µg/L	Result µg/L	COMPOUND	Quantitation Limit µg/L	Result µg/L
Benzene	2	U	1,3-Dichloropropane	2	U
Bromobenzene	2	U	cis-1,3-Dichloropropene	2	U
Bromochloromethane	2	U	trans-1,3-Dichloropropene	2	U
Bromodichloromethane	2	U	2,2-Dichloropropane	2	U
Bromoform	2	U	1,1-Dichloropropene	2	U
Bromomethane	2	U	Ethylbenzene	2	U
n-butylbenzene	2	U	Hexachlorobutadiene	2	U
sec-butylbenzene	2	U	Isopropylbenzene	2	U
tert-butylbenzene	2	U	p-isopropyltoluene	2	U
Carbon Tetrachloride	2	U	Methylene Chloride	5	U
Chlorobenzene	2	U	Methyl-tert-butyl ether (MTBE)	2	U
Chloroethane	2	U	Naphthalene	2	U
Chloroform	2	U	n-Propylbenzene	2	U
Chloromethane	2	U	Styrene	2	U
2-Chlorotoluene	2	U	1,1,1,2-Tetrachloroethane	2	U
4-Chlorotoluene	2	U	1,1,2,2-Tetrachloroethane	2	U
Dibromochloromethane	2	U	Tetrachloroethene	2	U
1,2-Dibromo-3-chloropropane	2	U	Toluene	2	U
1,2-Dibromoethane	2	U	1,2,3-Trichlorobenzene	2	U
Dibromomethane	2	U	1,2,4-Trichlorobenzene	2	U
1,2-Dichlorobenzene	2	U	1,1,1-Trichloroethane	2	U
1,3-Dichlorobenzene	2	U	1,1,2-Trichloroethane	2	U
1,4-Dichlorobenzene	2	U	Trichloroethene	2	U
Dichlorodifluoromethane	2	U	Trichlorofluoromethane	2	U
1,1-Dichloroethane	2	U	1,2,3-Trichloropropane	2	U
1,2-Dichloroethane	2	U	1,2,4-Trimethylbenzene	2	1 J
1,1-Dichloroethene	2	U	1,3,5-Trimethylbenzene	2	U
cis-1,2-Dichloroethene	2	U	Vinyl Chloride	2	U
trans-1,2-Dichloroethene	2	U	o-Xylene	2	U
1,2-Dichloropropane	2	U	m,p-Xylene	2	U
Acetone	10	U	Diethyl ether	2	U
Carbon Disulfide	2	U	2-Hexanone	10	U
Tetrahydrofuran	5	U	Methyl isobutyl ketone	10	U
Methyl ethyl ketone	10	U	Di-isopropyl ether (DIPE)	2	U
t-Butyl alcohol (TBA)	20	U	Ethyl t-butyl ether (ETBE)	2	U
t-Amyl methyl ether (TAME)	2	U			

Surrogate Standard Recovery

d4-1,2-Dichloroethane	101	%	d8-Toluene	100	%	Bromofluorobenzene	99 %

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis was conducted according to: Test Methods for Evaluating Solid Waste, SW-846 Method 8260B.

COMMENTS:

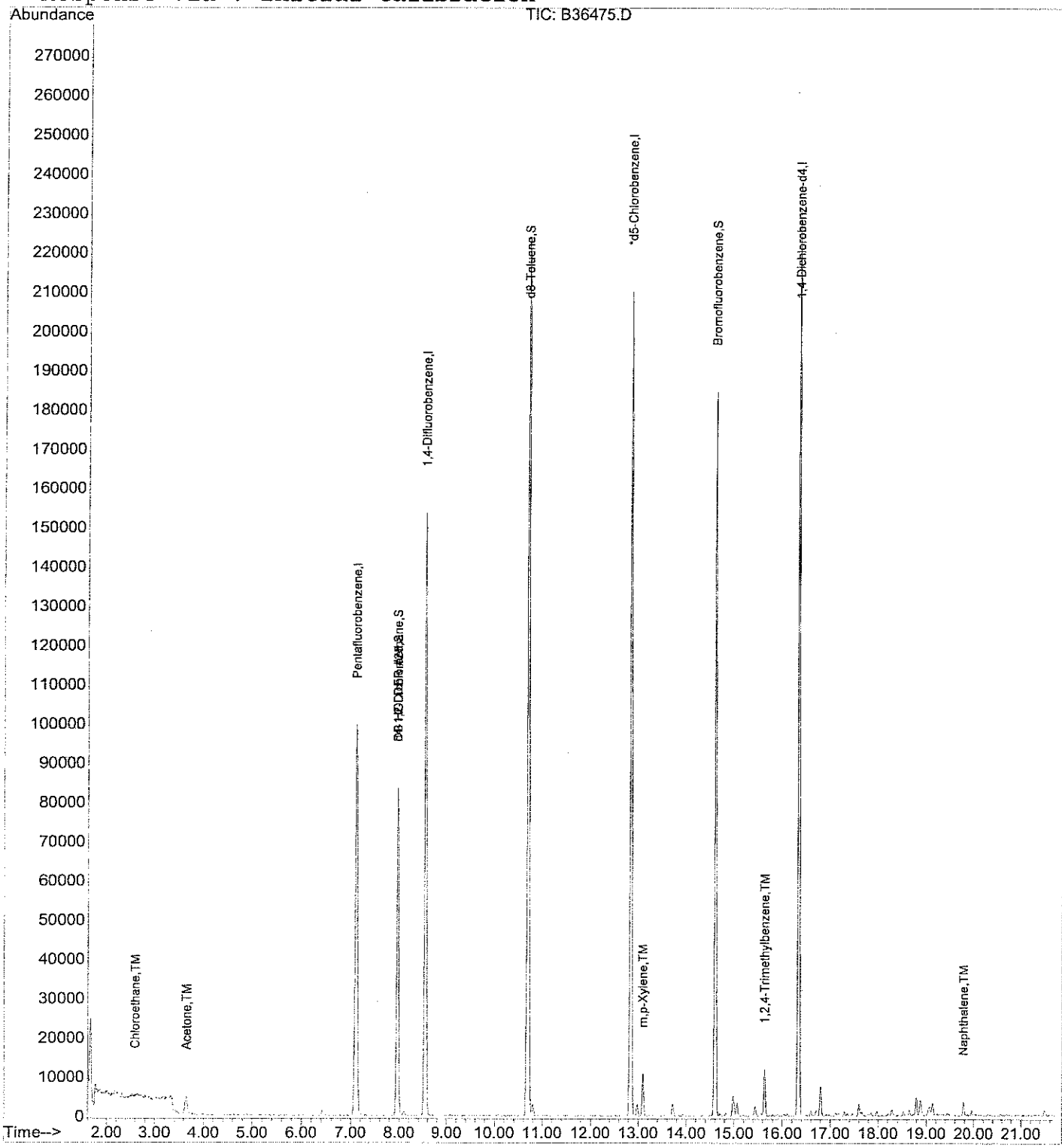
8260 full

Authorized signature

Quantitation Report

Data File : C:\HPCHEM\1\DATA\053006-B\B36475.D Vial: 19
Acq On : 30 May 2006 5:21 pm Operator:
Sample : 56637-20 Inst : Instrumen
Misc : 5000 Multiplr: 1.00
MS Integration Params: rteint.p
Quant Time: May 30 17:35 2006 Quant Results File: V805266B.RES

Method : C:\HPCHEM\1\METHODS\V805266B.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Tue May 30 08:22:32 2006
Response via : Initial Calibration



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May 31, 2006

CLIENT SAMPLE ID

Project Name: GP Old Town Mill

Project Number: 218736.002

Field Sample ID: MW-3

SAMPLE DATA

Lab Sample ID: 56637-21
Matrix: Aqueous
Percent Solid: N/A
Dilution Factor: 1
Collection Date: 05/26/06
Lab Receipt Date: 05/27/06
Analysis Date: 05/30/06

ANALYTICAL RESULTS VOLATILE ORGANICS

COMPOUND	Quantitation Limit µg/L	Result µg/L	COMPOUND	Quantitation Limit µg/L	Result µg/L
Benzene	2	U	1,3-Dichloropropane	2	U
Bromobenzene	2	U	cis-1,3-Dichloropropene	2	U
Bromochloromethane	2	U	trans-1,3-Dichloropropene	2	U
Bromodichloromethane	2	U	2,2-Dichloropropane	2	U
Bromoform	2	U	1,1-Dichloropropene	2	U
Bromomethane	2	U	Ethylbenzene	2	U
n-butylbenzene	2	U	Hexachlorobutadiene	2	U
sec-butylbenzene	2	U	Isopropylbenzene	2	U
tert-butylbenzene	2	U	p-isopropyltoluene	2	U
Carbon Tetrachloride	2	U	Methylene Chloride	5	U
Chlorobenzene	2	U	Methyl-tert-butyl ether (MTBE)	2	U
Chloroethane	2	U	Naphthalene	2	U
Chloroform	2	U	n-Propylbenzene	2	U
Chloromethane	2	U	Styrene	2	U
2-Chlorotoluene	2	U	1,1,1,2-Tetrachloroethane	2	U
4-Chlorotoluene	2	U	1,1,2,2-Tetrachloroethane	2	U
Dibromochloromethane	2	U	Tetrachloroethene	2	U
1,2-Dibromo-3-chloropropane	2	U	Toluene	2	U
1,2-Dibromoethane	2	U	1,2,3-Trichlorobenzene	2	U
Dibromomethane	2	U	1,2,4-Trichlorobenzene	2	U
1,2-Dichlorobenzene	2	U	1,1,1-Trichloroethane	2	U
1,3-Dichlorobenzene	2	U	1,1,2-Trichloroethane	2	U
1,4-Dichlorobenzene	2	U	Trichloroethene	2	U
Dichlorodifluoromethane	2	U	Trichlorofluoromethane	2	U
1,1-Dichloroethane	2	U	1,2,3-Trichloropropane	2	U
1,2-Dichloroethane	2	U	1,2,4-Trimethylbenzene	2	U
1,1-Dichloroethene	2	U	1,3,5-Trimethylbenzene	2	U
cis-1,2-Dichloroethene	2	U	Vinyl Chloride	2	U
trans-1,2-Dichloroethene	2	U	o-Xylene	2	U
1,2-Dichloropropane	2	U	m,p-Xylene	2	U
Acetone	10	U	Diethyl ether	2	U
Carbon Disulfide	2	U	2-Hexanone	10	U
Tetrahydrofuran	5	U	Methyl isobutyl ketone	10	U
Methyl ethyl ketone	10	U	Di-isopropyl ether (DIPE)	2	U
t-Butyl alcohol (TBA)	20	U	Ethyl t-butyl ether (ETBE)	2	U
t-Amyl methyl ether (TAME)	2	U			

Surrogate Standard Recovery

d4-1,2-Dichloroethane	101	%	d8-Toluene	100	%	Bromofluorobenzene	100 %

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis was conducted according to: Test Methods for Evaluating Solid Waste, SW-846 Method 8260B.

COMMENTS:

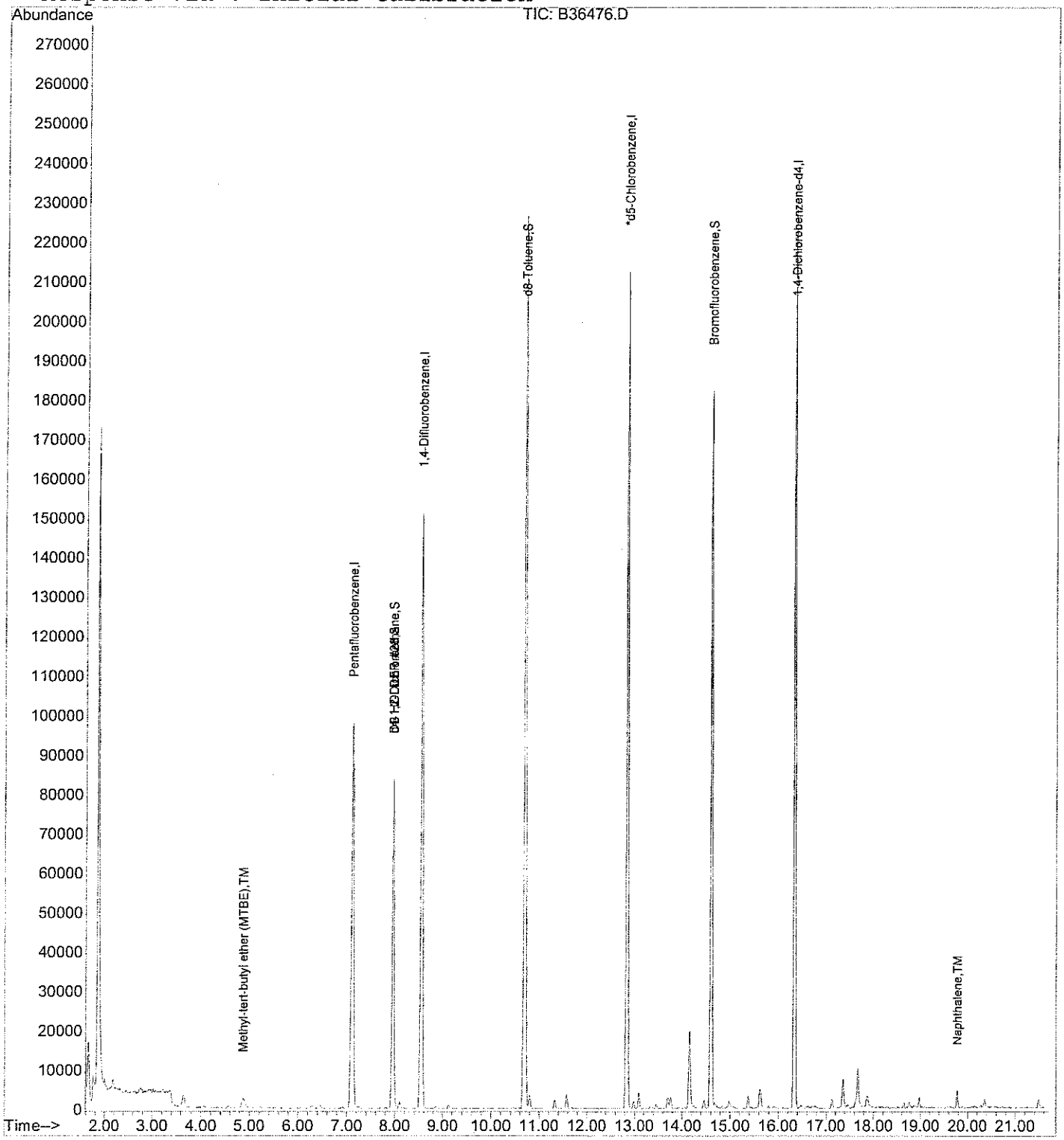
8260 full

Authorized signature Melissa Tull

Quantitation Report

Data File : C:\HPCHEM\1\DATA\053006-B\B36476.D Vial: 20
Acq On : 30 May 2006 5:50 pm Operator:
Sample : 56637-21 Inst : Instrumen
Misc : 5000 Multiplr: 1.00
MS Integration Params: rteint.p
Quant Time: May 30 18:05 2006 Quant Results File: V805266B.RES

Method : C:\HPCHEM\1\METHODS\V805266B.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Tue May 30 08:22:32 2006
Response via : Initial Calibration



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May 31, 2006

CLIENT SAMPLE ID

Project Name: GP Old Town Mill

Project Number: 218736.002

Field Sample ID: SB-40 (1 1/2-5 1/2)

SAMPLE DATA

Lab Sample ID: 56637-24
Matrix: Aqueous
Percent Solid: N/A
Dilution Factor: 1
Collection Date: 05/26/06
Lab Receipt Date: 05/27/06
Analysis Date: 05/30/06

ANALYTICAL RESULTS VOLATILE ORGANICS

COMPOUND	Quantitation Limit µg/L	Result µg/L	COMPOUND	Quantitation Limit µg/L	Result µg/L
Benzene	2	U	1,3-Dichloropropane	2	U
Bromobenzene	2	U	cis-1,3-Dichloropropene	2	U
Bromochloromethane	2	U	trans-1,3-Dichloropropene	2	U
Bromodichloromethane	2	U	2,2-Dichloropropane	2	U
Bromoform	2	U	1,1-Dichloropropene	2	U
Bromomethane	2	U	Ethylbenzene	2	U
n-butylbenzene	2	U	Hexachlorobutadiene	2	U
sec-butylbenzene	2	U	Isopropylbenzene	2	U
tert-butylbenzene	2	U	p-isopropyltoluene	2	U
Carbon Tetrachloride	2	U	Methylene Chloride	5	U
Chlorobenzene	2	U	Methyl-tert-butyl ether (MTBE)	2	U
Chloroethane	2	U	Naphthalene	2	U
Chloroform	2	U	n-Propylbenzene	2	U
Chloromethane	2	U	Styrene	2	U
2-Chlorotoluene	2	U	1,1,1,2-Tetrachloroethane	2	U
4-Chlorotoluene	2	U	1,1,2,2-Tetrachloroethane	2	U
Dibromochloromethane	2	U	Tetrachloroethene	2	U
1,2-Dibromo-3-chloropropane	2	U	Toluene	2	U
1,2-Dibromoethane	2	U	1,2,3-Trichlorobenzene	2	U
Dibromomethane	2	U	1,2,4-Trichlorobenzene	2	U
1,2-Dichlorobenzene	2	U	1,1,1-Trichloroethane	2	U
1,3-Dichlorobenzene	2	U	1,1,2-Trichloroethane	2	U
1,4-Dichlorobenzene	2	U	Trichloroethene	2	U
Dichlorodifluoromethane	2	U	Trichlorofluoromethane	2	U
1,1-Dichloroethane	2	U	1,2,3-Trichloropropane	2	U
1,2-Dichloroethane	2	U	1,2,4-Trimethylbenzene	2	U
1,1-Dichloroethene	2	U	1,3,5-Trimethylbenzene	2	U
cis-1,2-Dichloroethene	2	U	Vinyl Chloride	2	U
trans-1,2-Dichloroethene	2	U	o-Xylene	2	U
1,2-Dichloropropane	2	U	m,p-Xylene	2	U
Acetone	10	U	Diethyl ether	2	U
Carbon Disulfide	2	U	2-Hexanone	10	U
Tetrahydrofuran	5	U	Methyl isobutyl ketone	10	U
Methyl ethyl ketone	10	U	Di-isopropyl ether (DIPE)	2	U
t-Butyl alcohol (TBA)	20	U	Ethyl t-butyl ether (ETBE)	2	U
t-Amyl methyl ether (TAME)	2	U			

Surrogate Standard Recovery

d4-1,2-Dichloroethane	102 %	d8-Toluene	99 %	Bromofluorobenzene	101 %

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis was conducted according to: Test Methods for Evaluating Solid Waste, SW-846 Method 8260B.

COMMENTS:

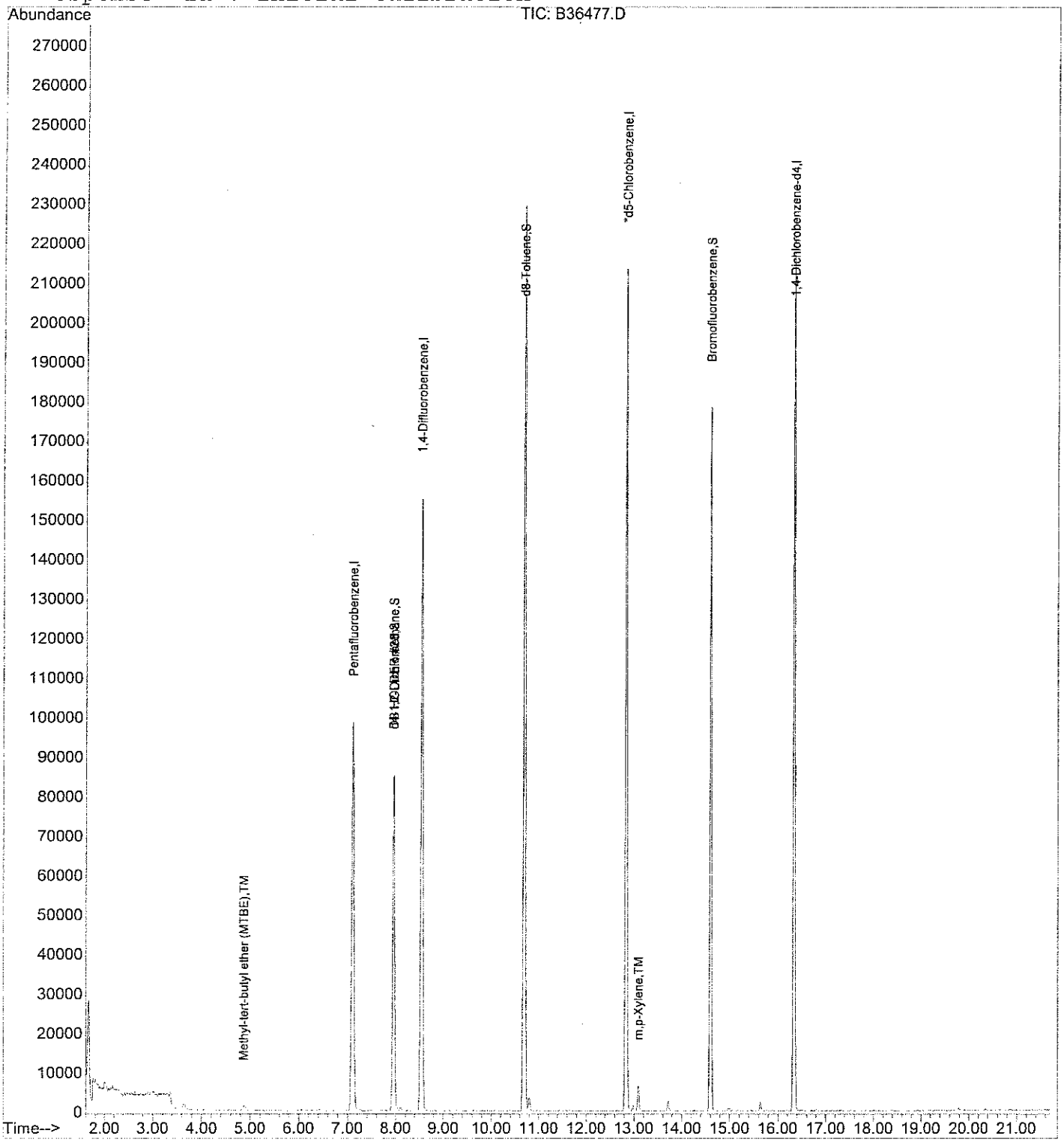
8260 full

Authorized signature

Quantitation Report

Data File : C:\HPCHEM\1\DATA\053006-B\B36477.D Vial: 21
Acq On : 30 May 2006 6:19 pm Operator:
Sample : 56637-24 Inst : Instrumen
Misc : 5000 Multiplr: 1.00
MS Integration Params: rteint.p
Quant Time: May 30 18:34 2006 Quant Results File: V805266B.RES

Method : C:\HPCHEM\1\METHODS\V805266B.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Tue May 30 08:22:32 2006
Response via : Initial Calibration



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May 31, 2006

CLIENT SAMPLE ID

Project Name: GP Old Town Mill

Project Number: 218736.002

Field Sample ID: SB-38 DUP

SAMPLE DATA

Lab Sample ID: 56637-25
Matrix: Aqueous
Percent Solid: N/A
Dilution Factor: 1
Collection Date: 05/25/06
Lab Receipt Date: 05/27/06
Analysis Date: 05/30/06

ANALYTICAL RESULTS VOLATILE ORGANICS

COMPOUND	Quantitation Limit µg/L	Result µg/L	COMPOUND	Quantitation Limit µg/L	Result µg/L
Benzene	2	U	1,3-Dichloropropane	2	U
Bromobenzene	2	U	cis-1,3-Dichloropropene	2	U
Bromochloromethane	2	U	trans-1,3-Dichloropropene	2	U
Bromodichloromethane	2	U	2,2-Dichloropropane	2	U
Bromoform	2	U	1,1-Dichloropropene	2	U
Bromomethane	2	U	Ethylbenzene	2	U
n-butylbenzene	2	U	Hexachlorobutadiene	2	U
sec-butylbenzene	2	U	Isopropylbenzene	2	U
tert-butylbenzene	2	U	p-isopropyltoluene	2	U
Carbon Tetrachloride	2	U	Methylene Chloride	5	U
Chlorobenzene	2	U	Methyl-tert-butyl ether (MTBE)	2	U
Chloroethane	2	U	Naphthalene	2	5
Chloroform	2	U	n-Propylbenzene	2	U
Chloromethane	2	U	Styrene	2	U
2-Chlorotoluene	2	U	1,1,1,2-Tetrachloroethane	2	U
4-Chlorotoluene	2	U	1,1,2,2-Tetrachloroethane	2	U
Dibromochloromethane	2	U	Tetrachloroethene	2	U
1,2-Dibromo-3-chloropropane	2	U	Toluene	2	U
1,2-Dibromoethane	2	U	1,2,3-Trichlorobenzene	2	U
Dibromomethane	2	U	1,2,4-Trichlorobenzene	2	U
1,2-Dichlorobenzene	2	U	1,1,1-Trichloroethane	2	U
1,3-Dichlorobenzene	2	U	1,1,2-Trichloroethane	2	U
1,4-Dichlorobenzene	2	U	Trichloroethene	2	U
Dichlorodifluoromethane	2	U	Trichlorofluoromethane	2	U
1,1-Dichloroethane	2	U	1,2,3-Trichloropropane	2	U
1,2-Dichloroethane	2	U	1,2,4-Trimethylbenzene	2	U
1,1-Dichloroethene	2	U	1,3,5-Trimethylbenzene	2	U
cis-1,2-Dichloroethene	2	U	Vinyl Chloride	2	U
trans-1,2-Dichloroethene	2	U	o-Xylene	2	U
1,2-Dichloropropane	2	U	m,p-Xylene	2	U
Acetone	10	U	Diethyl ether	2	U
Carbon Disulfide	2	U	2-Hexanone	10	U
Tetrahydrofuran	5	U	Methyl isobutyl ketone	10	U
Methyl ethyl ketone	10	U	Di-isopropyl ether (DIPE)	2	U
t-Butyl alcohol (TBA)	20	U	Ethyl t-butyl ether (ETBE)	2	U
t-Amyl methyl ether (TAME)	2	U			

Surrogate Standard Recovery

d4-1,2-Dichloroethane	101	%	d8-Toluene	99	%	Bromofluorobenzene	101	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis was conducted according to: Test Methods for Evaluating Solid Waste, SW-846 Method 8260B.

COMMENTS:

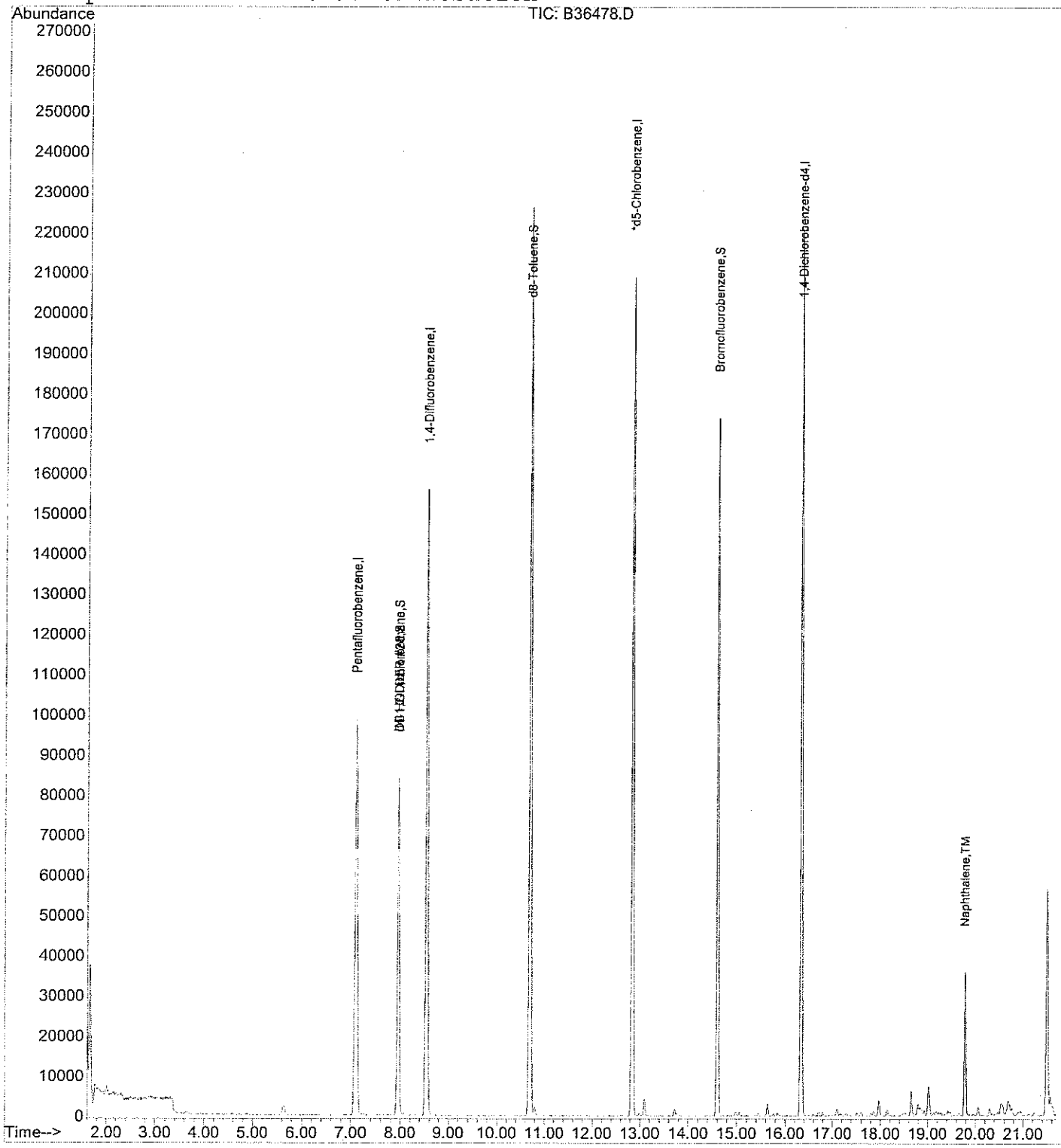
8260 full

Authorized signature

Quantitation Report

Data File : C:\HPCHEM\1\DATA\053006-B\B36478.D Vial: 22
Acq On : 30 May 2006 6:48 pm Operator:
Sample : 56637-25 Inst : Instrumen
Misc : 5000 Multiplr: 1.00
MS Integration Params: rteint.p
Quant Time: May 30 19:03 2006 Quant Results File: V805266B.RES

Method : C:\HPCHEM\1\METHODS\V805266B.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Tue May 30 08:22:32 2006
Response via : Initial Calibration



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May 31, 2006

CLIENT SAMPLE ID

Project Name: GP Old Town Mill

Project Number: 218736.002

Field Sample ID: MW-5

SAMPLE DATA

Lab Sample ID: 56637-26
Matrix: Aqueous
Percent Solid: N/A
Dilution Factor: 1
Collection Date: 05/26/06
Lab Receipt Date: 05/27/06
Analysis Date: 05/30/06

ANALYTICAL RESULTS VOLATILE ORGANICS

COMPOUND	Quantitation Limit µg/L	Result µg/L	COMPOUND	Quantitation Limit µg/L	Result µg/L
Benzene	2	U	1,3-Dichloropropane	2	U
Bromobenzene	2	U	cis-1,3-Dichloropropene	2	U
Bromochloromethane	2	U	trans-1,3-Dichloropropene	2	U
Bromodichloromethane	2	U	2,2-Dichloropropane	2	U
Bromoform	2	U	1,1-Dichloropropene	2	U
Bromomethane	2	U	Ethylbenzene	2	U
n-butylbenzene	2	U	Hexachlorobutadiene	2	U
sec-butylbenzene	2	U	Isopropylbenzene	2	U
tert-butylbenzene	2	U	p-isopropyltoluene	2	U
Carbon Tetrachloride	2	U	Methylene Chloride	5	U
Chlorobenzene	2	U	Methyl-tert-butyl ether (MTBE)	2	U
Chloroethane	2	U	Naphthalene	2	U
Chloroform	2	U	n-Propylbenzene	2	U
Chloromethane	2	U	Styrene	2	U
2-Chlorotoluene	2	U	1,1,1,2-Tetrachloroethane	2	U
4-Chlorotoluene	2	U	1,1,2,2-Tetrachloroethane	2	U
Dibromochloromethane	2	U	Tetrachloroethene	2	U
1,2-Dibromo-3-chloropropane	2	U	Toluene	2	U
1,2-Dibromoethane	2	U	1,2,3-Trichlorobenzene	2	U
Dibromomethane	2	U	1,2,4-Trichlorobenzene	2	U
1,2-Dichlorobenzene	2	U	1,1,1-Trichloroethane	2	U
1,3-Dichlorobenzene	2	U	1,1,2-Trichloroethane	2	U
1,4-Dichlorobenzene	2	U	Trichloroethene	2	U
Dichlorodifluoromethane	2	U	Trichlorofluoromethane	2	U
1,1-Dichloroethane	2	U	1,2,3-Trichloropropane	2	U
1,2-Dichloroethane	2	U	1,2,4-Trimethylbenzene	2	U
1,1-Dichloroethene	2	U	1,3,5-Trimethylbenzene	2	U
cis-1,2-Dichloroethene	2	U	Vinyl Chloride	2	U
trans-1,2-Dichloroethene	2	U	o-Xylene	2	U
1,2-Dichloropropane	2	U	m,p-Xylene	2	U
Acetone	10	U	Diethyl ether	2	U
Carbon Disulfide	2	U	2-Hexanone	10	U
Tetrahydrofuran	5	U	Methyl isobutyl ketone	10	U
Methyl ethyl ketone	10	U	Di-isopropyl ether (DIPE)	2	U
t-Butyl alcohol (TBA)	20	U	Ethyl t-butyl ether (ETBE)	2	U
t-Amyl methyl ether (TAME)	2	U			

Surrogate Standard Recovery

d4-1,2-Dichloroethane	102	%	d8-Toluene	100	%	Bromofluorobenzene	100	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis was conducted according to: Test Methods for Evaluating Solid Waste, SW-846 Method 8260B.

COMMENTS:

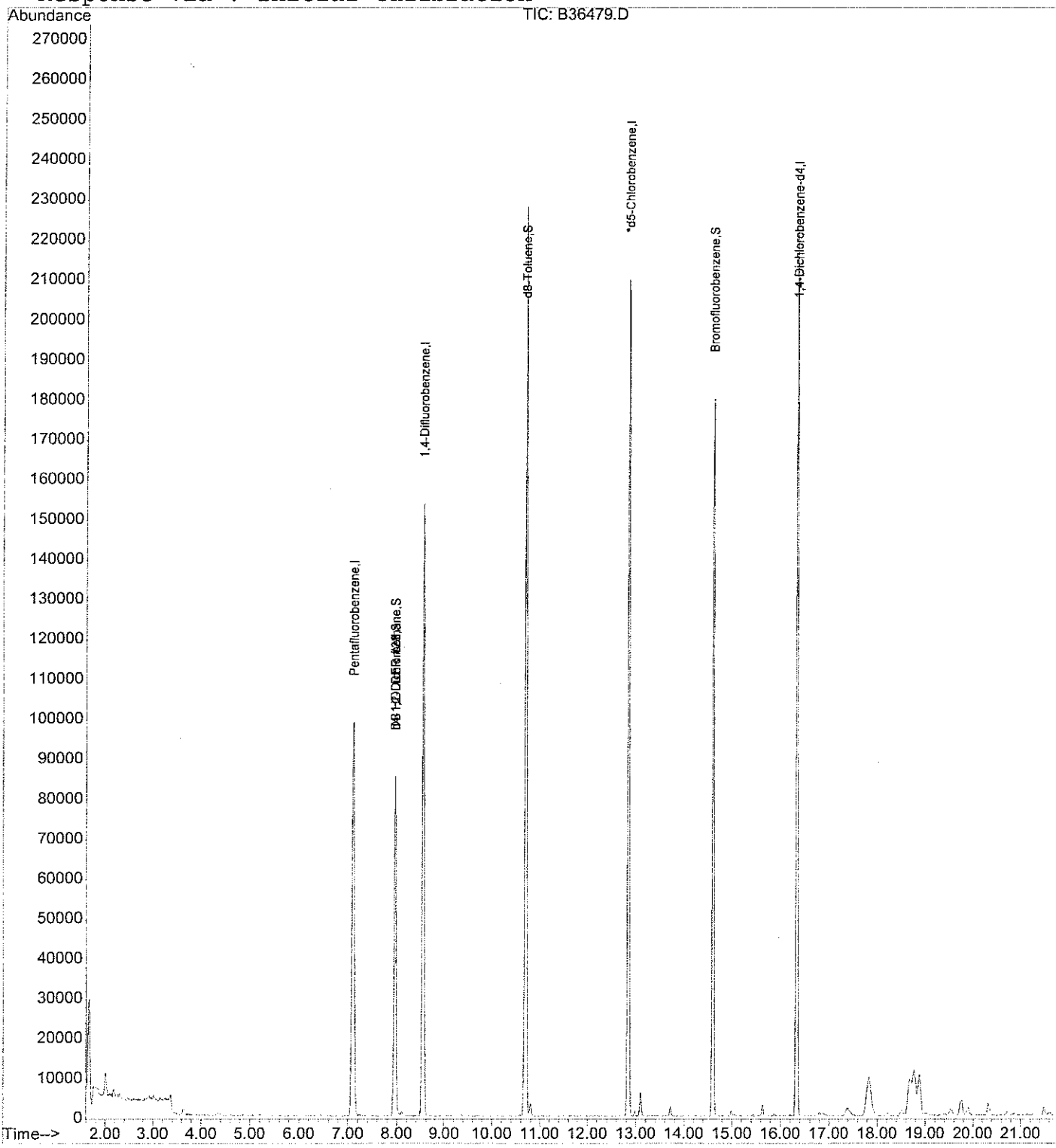
8260 full

Authorized signature

Quantitation Report

Data File : C:\HPCHEM\1\DATA\053006-B\B36479.D Vial: 23
Acq On : 30 May 2006 7:18 pm Operator:
Sample : 56637-26 Inst : Instrumen
Misc : 5000 Multiplr: 1.00
MS Integration Params: rteint.p
Quant Time: May 30 19:32 2006 Quant Results File: V805266B.RES

Method : C:\HPCHEM\1\METHODS\V805266B.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Tue May 30 08:22:32 2006
Response via : Initial Calibration



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May 31, 2006

CLIENT SAMPLE ID

Project Name: GP Old Town Mill

Project Number: 218736.002

Field Sample ID: SB-30

SAMPLE DATA

Lab Sample ID: 56637-27
Matrix: Aqueous
Percent Solid: N/A
Dilution Factor: 1
Collection Date: 05/25/06
Lab Receipt Date: 05/27/06
Analysis Date: 05/30/06

ANALYTICAL RESULTS VOLATILE ORGANICS

COMPOUND	Quantitation Limit µg/L	Result µg/L	COMPOUND	Quantitation Limit µg/L	Result µg/L
Benzene	2	U	1,3-Dichloropropane	2	U
Bromobenzene	2	U	cis-1,3-Dichloropropene	2	U
Bromochloromethane	2	U	trans-1,3-Dichloropropene	2	U
Bromodichloromethane	2	U	2,2-Dichloropropane	2	U
Bromoform	2	U	1,1-Dichloropropene	2	U
Bromomethane	2	U	Ethylbenzene	2	U
n-butylbenzene	2	U	Hexachlorobutadiene	2	U
sec-butylbenzene	2	U	Isopropylbenzene	2	U
tert-butylbenzene	2	U	p-isopropyltoluene	2	U
Carbon Tetrachloride	2	U	Methylene Chloride	5	U
Chlorobenzene	2	U	Methyl-tert-butyl ether (MTBE)	2	U
Chloroethane	2	U	Naphthalene	2	U
Chloroform	2	U	n-Propylbenzene	2	U
Chloromethane	2	U	Styrene	2	U
2-Chlorotoluene	2	U	1,1,1,2-Tetrachloroethane	2	U
4-Chlorotoluene	2	U	1,1,2,2-Tetrachloroethane	2	U
Dibromochloromethane	2	U	Tetrachloroethene	2	U
1,2-Dibromo-3-chloropropane	2	U	Toluene	2	U
1,2-Dibromoethane	2	U	1,2,3-Trichlorobenzene	2	U
Dibromomethane	2	U	1,2,4-Trichlorobenzene	2	U
1,2-Dichlorobenzene	2	U	1,1,1-Trichloroethane	2	U
1,3-Dichlorobenzene	2	U	1,1,2-Trichloroethane	2	U
1,4-Dichlorobenzene	2	U	Trichloroethene	2	U
Dichlorodifluoromethane	2	U	Trichlorofluoromethane	2	U
1,1-Dichloroethane	2	U	1,2,3-Trichloropropane	2	U
1,2-Dichloroethane	2	U	1,2,4-Trimethylbenzene	2	U
1,1-Dichloroethene	2	U	1,3,5-Trimethylbenzene	2	U
cis-1,2-Dichloroethene	2	U	Vinyl Chloride	2	U
trans-1,2-Dichloroethene	2	U	o-Xylene	2	U
1,2-Dichloropropane	2	U	m,p-Xylene	2	U
Acetone	10	U	Diethyl ether	2	U
Carbon Disulfide	2	U	2-Hexanone	10	U
Tetrahydrofuran	5	U	Methyl isobutyl ketone	10	U
Methyl ethyl ketone	10	U	Di-isopropyl ether (DIPE)	2	U
t-Butyl alcohol (TBA)	20	U	Ethyl t-butyl ether (ETBE)	2	U
t-Amyl methyl ether (TAME)	2	U			

Surrogate Standard Recovery

d4-1,2-Dichloroethane	102	%	d8-Toluene	99	%	Bromofluorobenzene	100 %

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis was conducted according to: Test Methods for Evaluating Solid Waste, SW-846 Method 8260B.

COMMENTS:

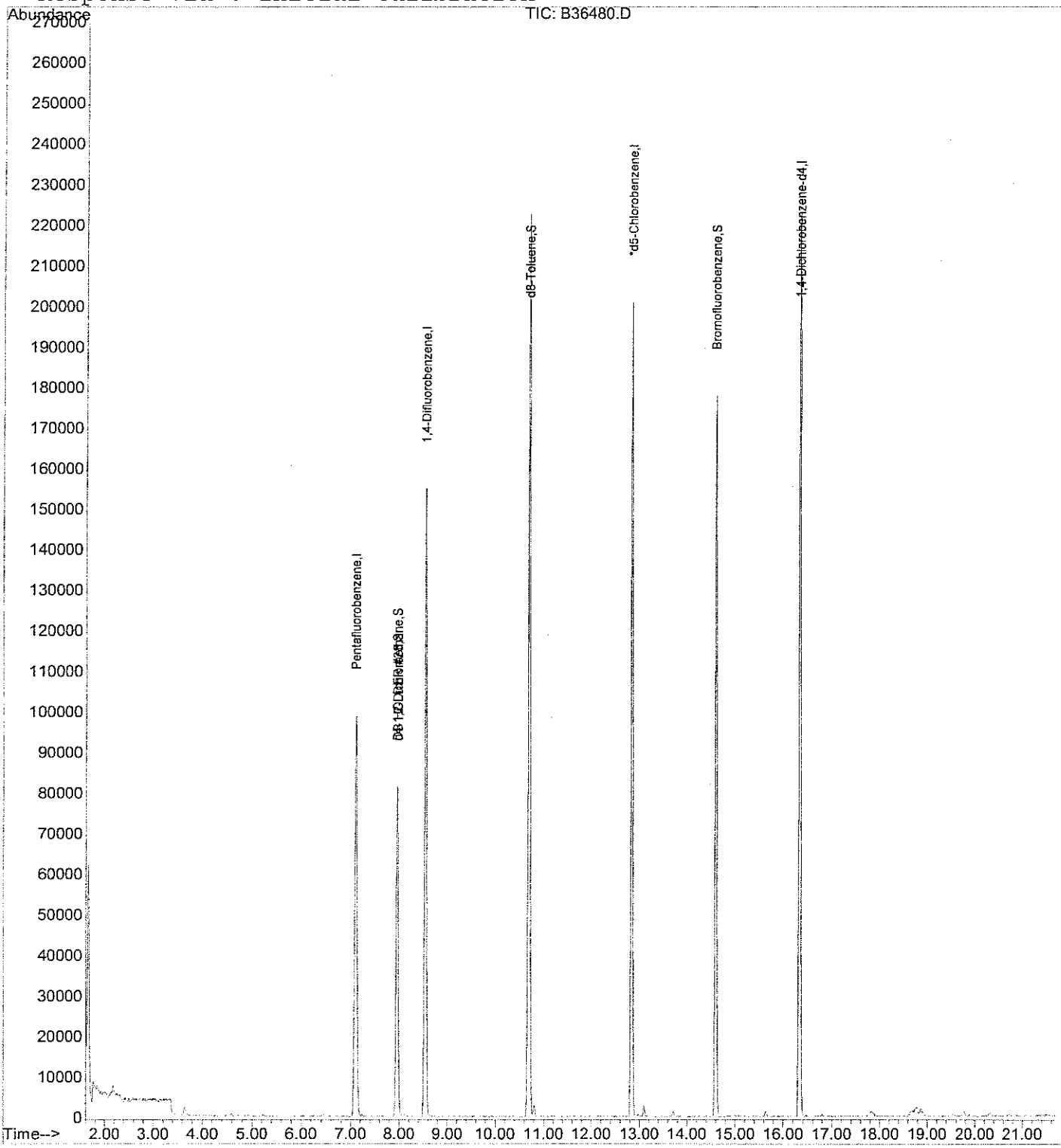
8260 full

Authorized signature

Quantitation Report

Data File : C:\HPCHEM\1\DATA\053006-B\B36480.D Vial: 24
Acq On : 30 May 2006 7:47 pm Operator:
Sample : 56637-27 Inst : Instrumen
Misc : 5000 Multiplr: 1.00
MS Integration Params: rteint.p
Quant Time: May 30 20:02 2006 Quant Results File: V805266B.RES

Method : C:\HPCHEM\1\METHODS\V805266B.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Tue May 30 08:22:32 2006
Response via : Initial Calibration



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May 31, 2006

CLIENT SAMPLE ID

Project Name: GP Old Town Mill

Project Number: 218736.002

Field Sample ID: SB-15

SAMPLE DATA

Lab Sample ID: 56637-28
Matrix: Aqueous
Percent Solid: N/A
Dilution Factor: 1
Collection Date: 05/25/06
Lab Receipt Date: 05/27/06
Analysis Date: 05/30/06

ANALYTICAL RESULTS VOLATILE ORGANICS

COMPOUND	Quantitation Limit µg/L	Result µg/L	COMPOUND	Quantitation Limit µg/L	Result µg/L
Benzene	2	U	1,3-Dichloropropane	2	U
Bromobenzene	2	U	cis-1,3-Dichloropropene	2	U
Bromochloromethane	2	U	trans-1,3-Dichloropropene	2	U
Bromodichloromethane	2	U	2,2-Dichloropropane	2	U
Bromoform	2	U	1,1-Dichloropropene	2	U
Bromomethane	2	U	Ethylbenzene	2	U
n-butylbenzene	2	U	Hexachlorobutadiene	2	U
sec-butylbenzene	2	U	Isopropylbenzene	2	U
tert-butylbenzene	2	U	p-isopropyltoluene	2	U
Carbon Tetrachloride	2	U	Methylene Chloride	5	U
Chlorobenzene	2	U	Methyl-tert-butyl ether (MTBE)	2	U
Chloroethane	2	U	Naphthalene	2	U
Chloroform	2	U	n-Propylbenzene	2	U
Chloromethane	2	U	Styrene	2	U
2-Chlorotoluene	2	U	1,1,1,2-Tetrachloroethane	2	U
4-Chlorotoluene	2	U	1,1,2,2-Tetrachloroethane	2	U
Dibromochloromethane	2	U	Tetrachloroethene	2	U
1,2-Dibromo-3-chloropropane	2	U	Toluene	2	U
1,2-Dibromoethane	2	U	1,2,3-Trichlorobenzene	2	U
Dibromomethane	2	U	1,2,4-Trichlorobenzene	2	U
1,2-Dichlorobenzene	2	U	1,1,1-Trichloroethane	2	U
1,3-Dichlorobenzene	2	U	1,1,2-Trichloroethane	2	U
1,4-Dichlorobenzene	2	U	Trichloroethene	2	U
Dichlorodifluoromethane	2	U	Trichlorofluoromethane	2	U
1,1-Dichloroethane	2	U	1,2,3-Trichloropropane	2	U
1,2-Dichloroethane	2	U	1,2,4-Trimethylbenzene	2	U
1,1-Dichloroethene	2	U	1,3,5-Trimethylbenzene	2	U
cis-1,2-Dichloroethene	2	U	Vinyl Chloride	2	U
trans-1,2-Dichloroethene	2	U	o-Xylene	2	U
1,2-Dichloropropane	2	U	m,p-Xylene	2	U
Acetone	10	**15**	Diethyl ether	2	U
Carbon Disulfide	2	U	2-Hexanone	10	U
Tetrahydrofuran	5	U	Methyl isobutyl ketone	10	U
Methyl ethyl ketone	10	U	Di-isopropyl ether (DIPE)	2	U
t-Butyl alcohol (TBA)	20	U	Ethyl t-butyl ether (ETBE)	2	U
t-Amyl methyl ether (TAME)	2	U			

Surrogate Standard Recovery

d4-1,2-Dichloroethane	102	%	d8-Toluene	101	%	Bromofluorobenzene	100 %

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis was conducted according to: Test Methods for Evaluating Solid Waste, SW-846 Method 8260B.

COMMENTS:

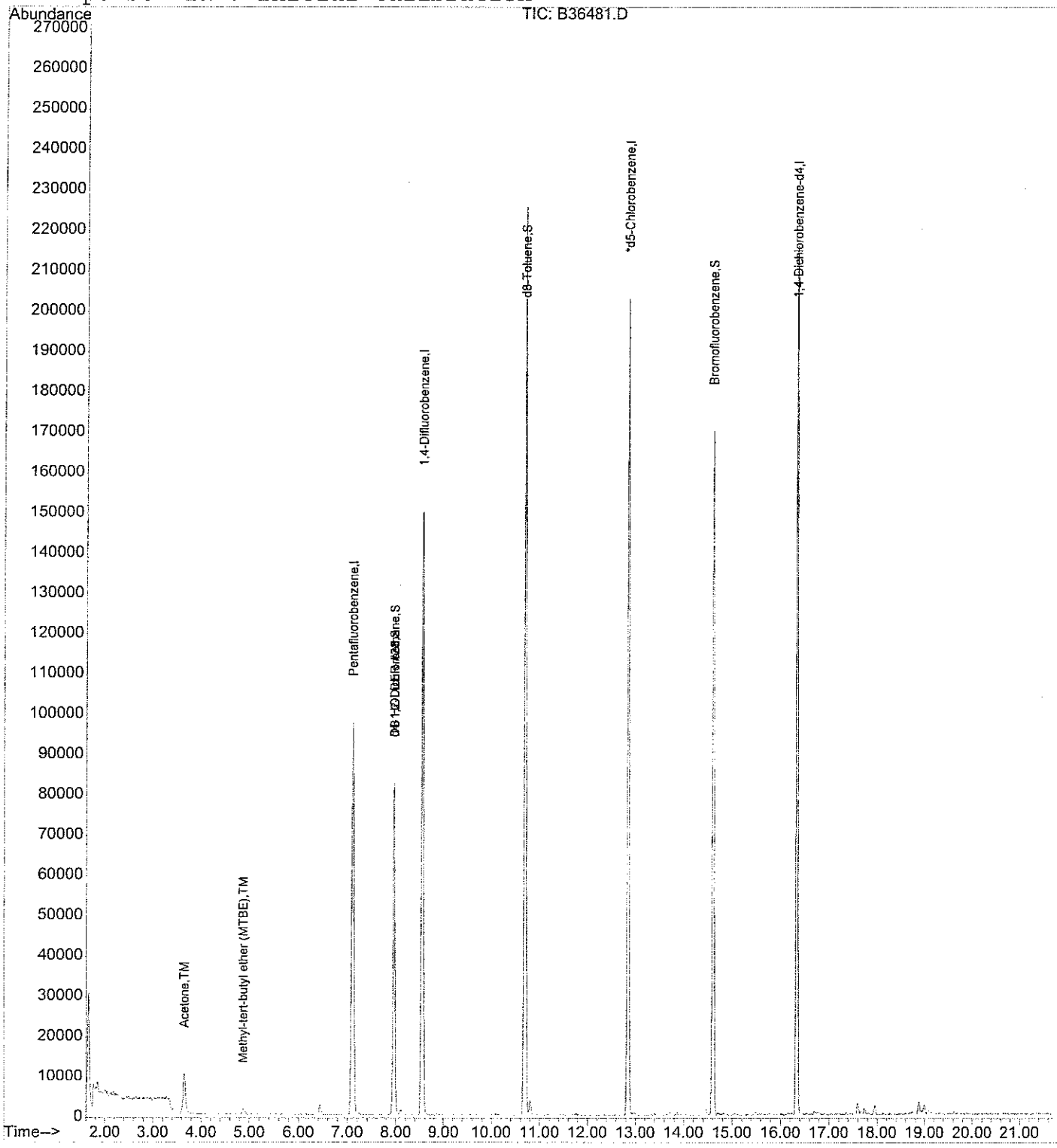
8260 full

Authorized signature

Quantitation Report

Data File : C:\HPCHEM\1\DATA\053006-B\B36481.D Vial: 25
Acq On : 30 May 2006 8:16 pm Operator:
Sample : 56637-28 Inst : Instrumen
Misc : 5000 Multiplr: 1.00
MS Integration Params: rteint.p
Quant Time: May 30 20:31 2006 Quant Results File: V805266B.RES

Method : C:\HPCHEM\1\METHODS\V805266B.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Tue May 30 08:22:32 2006
Response via : Initial Calibration



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May 31, 2006

CLIENT SAMPLE ID

Project Name: GP Old Town Mill

Project Number: 218736.002

Field Sample ID: Trip Blank

SAMPLE DATA

Lab Sample ID: 56637-29
Matrix: Aqueous
Percent Solid: N/A
Dilution Factor: 1
Collection Date: 05/26/06
Lab Receipt Date: 05/27/06
Analysis Date: 05/30/06

ANALYTICAL RESULTS VOLATILE ORGANICS

COMPOUND	Quantitation Limit µg/L	Result µg/L	COMPOUND	Quantitation Limit µg/L	Result µg/L
Benzene	2	U	1,3-Dichloropropane	2	U
Bromobenzene	2	U	cis-1,3-Dichloropropene	2	U
Bromochloromethane	2	U	trans-1,3-Dichloropropene	2	U
Bromodichloromethane	2	U	2,2-Dichloropropane	2	U
Bromoform	2	U	1,1-Dichloropropene	2	U
Bromomethane	2	U	Ethylbenzene	2	U
n-butylbenzene	2	U	Hexachlorobutadiene	2	U
sec-butylbenzene	2	U	Isopropylbenzene	2	U
tert-butylbenzene	2	U	p-isopropyltoluene	2	U
Carbon Tetrachloride	2	U	Methylene Chloride	5	U
Chlorobenzene	2	U	Methyl-tert-butyl ether (MTBE)	2	U
Chloroethane	2	U	Naphthalene	2	U
Chloroform	2	U	n-Propylbenzene	2	U
Chloromethane	2	U	Styrene	2	U
2-Chlorotoluene	2	U	1,1,1,2-Tetrachloroethane	2	U
4-Chlorotoluene	2	U	1,1,2,2-Tetrachloroethane	2	U
Dibromochloromethane	2	U	Tetrachloroethene	2	U
1,2-Dibromo-3-chloropropane	2	U	Toluene	2	U
1,2-Dibromoethane	2	U	1,2,3-Trichlorobenzene	2	U
Dibromomethane	2	U	1,2,4-Trichlorobenzene	2	U
1,2-Dichlorobenzene	2	U	1,1,1-Trichloroethane	2	U
1,3-Dichlorobenzene	2	U	1,1,2-Trichloroethane	2	U
1,4-Dichlorobenzene	2	U	Trichloroethene	2	U
Dichlorodifluoromethane	2	U	Trichlorofluoromethane	2	U
1,1-Dichloroethane	2	U	1,2,3-Trichloropropane	2	U
1,2-Dichloroethane	2	U	1,2,4-Trimethylbenzene	2	U
1,1-Dichloroethene	2	U	1,3,5-Trimethylbenzene	2	U
cis-1,2-Dichloroethene	2	U	Vinyl Chloride	2	U
trans-1,2-Dichloroethene	2	U	o-Xylene	2	U
1,2-Dichloropropane	2	U	m,p-Xylene	2	U
Acetone	10	U	Diethyl ether	2	U
Carbon Disulfide	2	U	2-Hexanone	10	U
Tetrahydrofuran	5	U	Methyl isobutyl ketone	10	U
Methyl ethyl ketone	10	U	Di-isopropyl ether (DIPE)	2	U
t-Butyl alcohol (TBA)	20	U	Ethyl t-butyl ether (ETBE)	2	U
t-Amyl methyl ether (TAME)	2	U			

Surrogate Standard Recovery

d4-1,2-Dichloroethane	100	%	d8-Toluene	99	%	Bromofluorobenzene	100	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis was conducted according to: Test Methods for Evaluating Solid Waste, SW-846 Method 8260B.

COMMENTS:

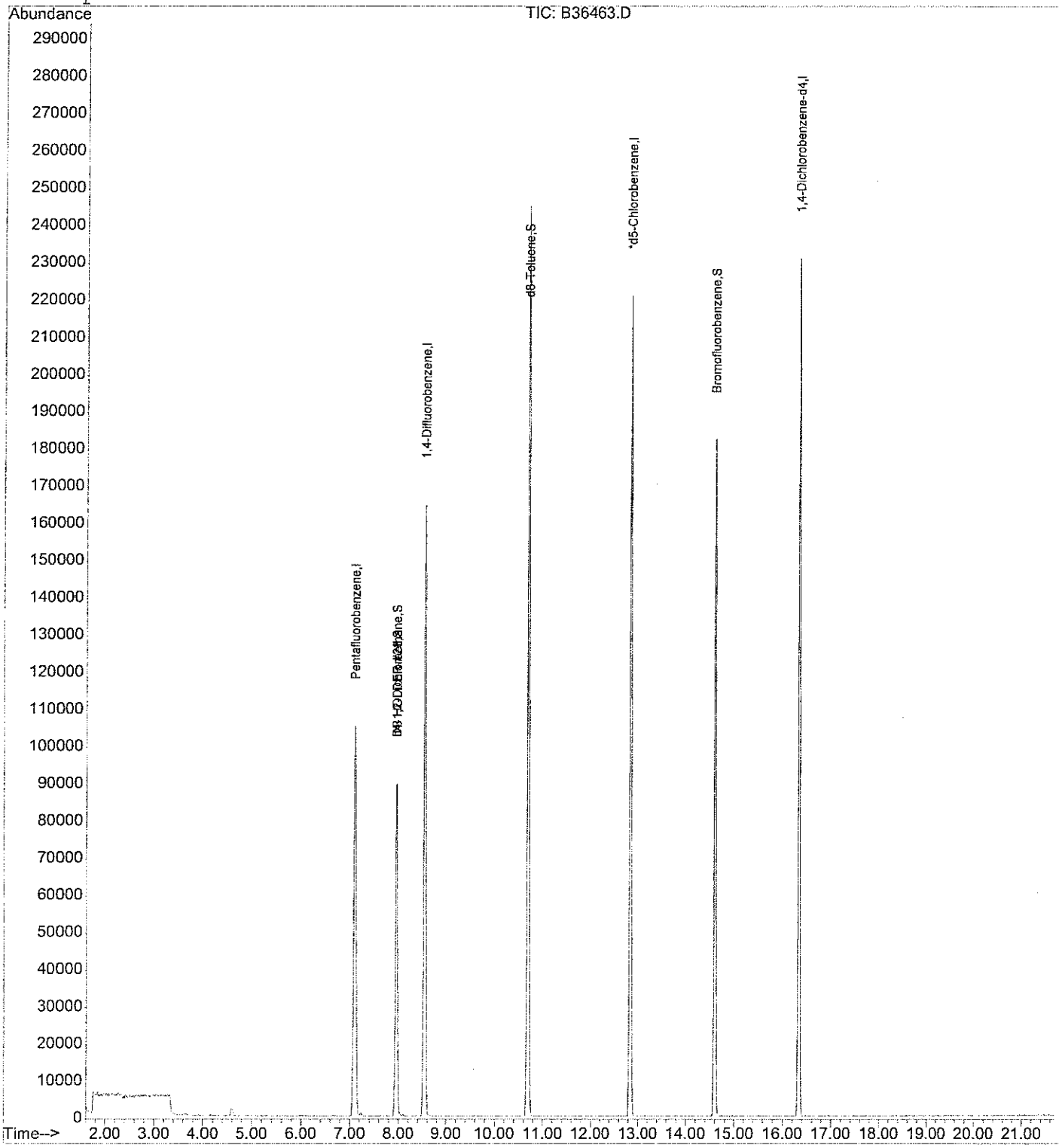
8260 full

Authorized signature

Quantitation Report

Data File : C:\HPCHEM\1\DATA\053006-B\B36463.D Vial: 7
Acq On : 30 May 2006 11:20 am Operator:
Sample : 56637-29 Inst : Instrumen
Misc : 5000 Multiplr: 1.00
MS Integration Params: rteint.p
Quant Time: May 30 11:41 2006 Quant Results File: V805266B.RES

Method : C:\HPCHEM\1\METHODS\V805266B.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Tue May 30 08:22:32 2006
Response via : Initial Calibration



analytics environmental laboratory LLC

VOLATILE
QC FORMS

analytics environmental laboratory LLC

VOLATILE ORGANIC SOLID
SYSTEM MONITORING COMPOUNDS
SUMMARY

Project #:	218736.002	Quote #:		SDG#:	56637
BFB File ID:	C59021T.D	Site ID:	GP Old Town Mill	Location:	ME
Instrument ID:	C			BFB Injection Date:	5/22/06
GC Column:	RTX-502.2	0.25(mm)		BFB Injection Time:	1030
				Heated Purge: (Y/N)	N

	SAMPLE ID	SMC 1 #	SMC 2 #	SMC 3 #	OUT
01	0.2 PPB	NA	NA	NA	0
02	1 PPB	NA	NA	NA	0
03	2 PPB	NA	NA	NA	0
04	5 PPB	NA	NA	NA	0
05	20 PPB	93	94	91	0
06	50 PPB	NA	NA	NA	0
07	100 PPB	NA	NA	NA	0
08	ZZZZZ	63 *	52 *	92	2
09	ZZZZZ	96	92	91	0
10	ZZZZZ	88	90	93	0
11	ZZZZZ	84	92	93	0
12	ZZZZZ	89	92	93	0
13	ZZZZZ	93	89	92	0
14	ZZZZZ	91	91	90	0
15	ZZZZZ	80	70	80	0
16	ZZZZZ	85	91	89	0
17	ZZZZZ	90	91	90	0
18	ZZZZZ	90	80	90	0
19	ZZZZZ	90	80	90	0
20	ZZZZZ	90	80	90	0
21	ZZZZZ	90	80	90	0
22	ZZZZZ	90	80	90	0
23	ZZZZZ	90	80	90	0
24	ZZZZZ	90	80	90	0
25					
25					
26					

	QC LIMIT
SMC1 = 1,2-Dichloroethane-d4	70-130%
SMC2 = Toluene-d8	70-130%
SMC3 = Bromofluorobenzene	70-130%

Column to be used to flag recovery values

* Values outside of protocol required QC limits

D System Monitoring Compound diluted out

analytics environmental laboratory LLC

VOLATILE ORGANIC AQUEOUS
SYSTEM MONITORING COMPOUNDS
SUMMARY

Project #:	218736.002	Quote #:		SDG#:	56637
BFB File ID:	B36407T.D	Site ID:	GP Old Town Mill	Location:	ME
Instrument ID:	B			BFB Injection Date:	5/26/06
GC Column:	RTX-502.2	ID: 0.25(mm)		BFB Injection Time:	0956
				Heated Purge: (Y/N)	N

	SAMPLE ID	SMC 1 #	SMC 2 #	SMC 3 #	OUT
01	1 PPB	103	100	101	0
02	2 PPB	103	99	99	0
03	5 PPB	102	99	100	0
04	20 PPB	101	100	101	0
05	50 PPB	95	101	100	0
06	100 PPB	96	100	99	0
07	ZZZZZ	100	99	99	0
08	ZZZZZ	99	99	100	0
09	ZZZZZ	98	100	101	0
10	ZZZZZ	100	99	99	0
11	ZZZZZ	102	99	99	0
12	ZZZZZ	102	99	98	0
13	ZZZZZ	102	99	100	0
14	ZZZZZ	103	100	99	0
15	ZZZZZ	103	99	99	0
16	ZZZZZ	102	99	98	0
17	ZZZZZ	101	99	101	0
18	ZZZZZ	99	100	102	0
19	ZZZZZ	102	99	99	0
20	ZZZZZ	103	99	99	0
21	ZZZZZ	102	101	97	0
22	ZZZZZ	102	100	99	0
23	ZZZZZ	102	99	98	0
24					
25					
25					
26					

	QC LIMIT
SMC1 = 1,2-Dichloroethane-d4	70-130%
SMC2 = Toluene-d8	70-130%
SMC3 = Bromofluorobenzene	70-130%

\# Column to be used to flag recovery values

* Values outside of protocol required QC limits

D System Monitoring Compound diluted out

analytics environmental laboratory LLC

VOLATILE ORGANIC SOLID
SYSTEM MONITORING COMPOUNDS
SUMMARY

Project #:	218736.002	Quote #:		SDG#:	56637
BFB File ID:	C59117T.D	Site ID:	GP Old Town Mill	Location:	ME
Instrument ID:	C			BFB Injection Date:	5/30/06
GC Column:	RTX-502.2	0.25(mm)		BFB Injection Time:	0818
				Heated Purge: (Y/N)	N

	SAMPLE ID	SMC 1 #	SMC 2 #	SMC 3 #	OUT
01	20 PPB	92	89	89	0
02	LS05306C	83	82	84	0
03	LS05306C2	75	82	82	0
04	ZZZZZ	89	91	93	0
05	MB05306C	76	84	84	0
06	56637-4	78	76	79	0
07	56637-14	64 *	69 *	67 *	3
08	56637-15	99	103	103	0
09	56637-16	78	79	84	0
10	ZZZZZ	91	86	88	0
11	56637-19	98	95	93	0
12	56637-9	98	95	93	0
13	56637-10	76	80	82	0
14	56637-11	71	78	79	0
15	ZZZZZ	62 *	68 *	70	0
16	ZZZZZ	75	66 *	69 *	2
17	56637-14, RR	60 *	63 *	58 *	3
18	56637-17, RR	73	78	82	0
19	ZZZZZ	63 *	68 *	72	2
20	56637-18, RR,, 10ML FV	71	70	72	0
21	ZZZZZ	59 *	63 *	63 *	3
22	ZZZZZ	49 *	38 *	38 *	3
23					
24					
25					
25					
26					

	QC LIMIT
SMC1 = 1,2-Dichloroethane-d4	70-130%
SMC2 = Toluene-d8	70-130%
SMC3 = Bromofluorobenzene	70-130%

Column to be used to flag recovery values

* Values outside of protocol required QC limits

D System Monitoring Compound diluted out

VOLATILE ORGANIC AQUEOUS
SYSTEM MONITORING COMPOUNDS
SUMMARY

Project #:	218736.002	Quote #:		SDG#:	56637
BFB File ID:	B36458T.D	Site ID:	GP Old Town Mill	Location:	ME
Instrument ID:	B			BFB Injection Date:	5/30/06
GC Column:	RTX-502.2	ID: 0.25(mm)		BFB Injection Time:	0858
				Heated Purge: (Y/N)	N

	SAMPLE ID	SMC 1 #	SMC 2 #	SMC 3 #	OUT
01	20 PPB	97	99	101	0
02	L805306B	96	100	102	0
03	ZZZZZ	96	100	98	0
04	B805306B	99	99	100	0
05	56637-29	100	99	100	0
06	ZZZZZ	99	100	101	0
07	ZZZZZ	100	99	99	0
08	ZZZZZ	103	100	98	0
09	ZZZZZ	103	100	100	0
10	ZZZZZ	101	100	99	0
11	56637-1	101	100	99	0
12	56637-2	102	100	102	0
13	56637-2,MS	100	101	101	0
14	56637-2,MSD	97	100	101	0
15	56637-3	102	99	99	0
16	56637-5	100	100	100	0
17	56637-20	101	100	99	0
18	56637-21	101	100	100	0
19	56637-24	102	99	101	0
20	56637-25	101	99	101	0
21	56637-26	102	100	100	0
22	56637-27	102	99	100	0
23	56637-28	102	101	100	0
24					
25					
25					
26					

	QC LIMIT
SMC1 = 1,2-Dichloroethane-d4	70-130%
SMC2 = Toluene-d8	70-130%
SMC3 = Bromofluorobenzene	70-130%

\# Column to be used to flag recovery values

* Values outside of protocol required QC limits

D System Monitoring Compound diluted out

analytics environmental laboratory LLC

VOLATILE ORGANIC

LABORATORY CONTROL SAMPLE

PERCENT RECOVERY

Project #: 218736.002 Quote #: SDG#: 56637

Site ID: GP Old Town Mill Location: ME

Instrument ID: B

GC Column: RTX-502.2 ID: 0.25 (mm)

Heated Purge: (Y/N) N

Laboratory Cotrol Sample ID: L805306B

COMPOUND	SPIKE ADDED (ug/L)	SAMPLE AMT (ug/L)	LCS CONC. (ug/L)	LCS REC %	#	QC Limits %REC
Dichlorodifluoromethane	20	U	23.0	115		80-120
Chloromethane	20	U	18.9	94		80-120
Vinyl Chloride	20	U	15.7	79	*	80-120
Bromomethane	20	U	20.2	101		80-120
Chloroethane	20	U	18.1	90		80-120
t-Butyl alcohol (TBA)	100	U	122.8	123		70-130
Trichlorofluoromethane	20	U	18.6	93		80-120
Diethyl ether	20	U	19.2	96		80-120
Acetone	100	U	99.8	100		70-130
1,1-Dichloroethene	20	U	19.0	95		80-120
Di-isopropyl ether (DIPE)	20	U	19.6	98		80-120
Methylene Chloride	20	U	20.0	100		80-120
Carbon Disulfide	20	U	17.3	86		80-120
Methyl-tert-butyl ether (MTBE)	40	U	41.5	104		80-120
trans-1,2-Dichloroethene	20	U	18.6	93		80-120
1,1-Dichloroethane	20	U	18.9	94		80-120
Methyl ethyl ketone	100	U	103.9	104		70-130
Ethyl t-butyl ether (ETBE)	20	U	19.8	99		80-120
2,2-Dichloropropane	20	U	22.0	110		80-120
cis-1,2-Dichloroethene	20	U	20.2	101		80-120
t-Amyl methyl ether (TAME)	20	U	20.5	103		80-120
Chloroform	20	U	19.2	96		80-120
Bromochloromethane	20	U	19.9	100		80-120
Tetrahydrofuran	20	U	21.1	105		70-130
1,1,1-Trichloroethane	20	U	20.1	100		80-120
1,1-Dichloropropene	20	U	19.4	97		80-120
Carbon Tetrachloride	20	U	21.5	107		80-120
1,2-Dichloroethane	20	U	19.0	95		80-120
Benzene	20	U	18.9	94		80-120
Trichloroethene	20	U	19.8	99		80-120
1,2-Dichloropropane	20	U	19.7	99		80-120
Bromodichloromethane	20	U	21.8	109		80-120
Dibromomethane	20	U	20.6	103		80-120
2-Hexanone	100	U	108.2	108		70-130
Methyl isobutyl ketone	100	U	104.8	105		70-130
cis-1,3-Dichloropropene	20	U	21.3	107		80-120
Toluene	20	U	19.2	96		80-120
trans-1,3-Dichloropropene	20	U	20.5	102		80-120
1,1,2-Trichloroethane	20	U	20.9	105		80-120
1,3-Dichloropropane	20	U	20.5	103		80-120
Tetrachloroethene	20	U	20.0	100		80-120
Dibromochloromethane	20	U	23.0	115		80-120
1,2-Dibromoethane	20	U	21.0	105		80-120
Chlorobenzene	20	U	19.5	97		80-120
1,1,1,2-Tetrachloroethane	20	U	22.2	111		80-120
Ethylbenzene	20	U	19.2	96		80-120
m,p-Xylene	40	U	38.6	96		80-120

analytics environmental laboratory LLC

VOLATILE ORGANIC

LABORATORY CONTROL SAMPLE

PERCENT RECOVERY

Project #: 218736.002 Quote #: SDG#: 56637

Site ID: GP Old Town Mill Location: ME

Instrument ID: B

GC Column: RTX-502.2 ID: 0.25 (mm)

Heated Purge: (Y/N) N

Laboratory Cotrol Sample ID: L805306B

COMPOUND	SPIKE ADDED (ug/L)	SAMPLE AMT (ug/L)	LCS CONC. (ug/L)	LCS REC %	#	QC Limits %REC
o-Xylene	20	U	19.9	99		80-120
Styrene	20	U	19.8	99		80-120
Bromoform	20	U	23.5	117		80-120
Isopropylbenzene	20	U	18.3	92		80-120
1,1,2,2-Tetrachloroethane	20	U	20.7	104		80-120
1,2,3-Trichloropropane	20	U	20.9	104		80-120
n-Propylbenzene	20	U	19.2	96		80-120
Bromobenzene	20	U	19.7	98		80-120
1,3,5-Trimethylbenzene	20	U	19.2	96		80-120
2-Chlorotoluene	20	U	19.5	97		80-120
4-Chlorotoluene	20	U	19.7	98		80-120
tert-butylbenzene	20	U	19.5	97		80-120
1,2,4-Trimethylbenzene	20	U	19.2	96		80-120
sec-butylbenzene	20	U	19.9	100		80-120
p-isopropyltoluene	20	U	19.2	96		80-120
1,3-Dichlorobenzene	20	U	19.6	98		80-120
1,4-Dichlorobenzene	20	U	19.4	97		80-120
n-butylbenzene	20	U	19.5	97		80-120
1,2-Dichlorobenzene	20	U	19.6	98		80-120
1,2-Dibromo-3-chloropropane	20	U	24.1	121	*	80-120
1,2,4-Trichlorobenzene	20	U	20.3	101		80-120
Hexachlorobutadiene	20	U	22.0	110		80-120
Naphthalene	20	U	21.5	107		80-120
1,2,3-Trichlorobenzene	20	U	20.6	103		80-120

Column to be used to flag recovery and RPD values with an asterisk

* Values outside of QC limits

Spike Recovery: 2 out of 71 outside limits

Comments: _____

_____

analytics environmental laboratory LLC

VOLATILE ORGANIC
MATRIX SPIKE/DUPLICATE
PERCENT RECOVERY

Project #: 218736.002 Quote #: SDG#: 56637

Site ID: GP Old Town Mill Location: ME

Instrument ID: B
GC Column: RTX-502.2 ID: 0.25 (mm)

Heated Purge: (Y/N) N

Matrix Spike/Duplicate Sample ID: 56637-2

COMPOUND	SPIKE ADDED (ug/L)	SAMPLE AMT (ug/L)	MS CONC. (ug/L)	MS REC %	#	MSD CONC. (ug/L)	MSD REC %	#	RPD	#	QC LIMITS RPD	QC LIMITS % REC.
Dichlorodifluoromethane	20	U	23.5	118		23.9	119		1		15	70-130
Chloromethane	20	U	18.7	93		17.8	89		4		15	70-130
Vinyl Chloride	20	U	17.2	86		17.9	89		4		15	70-130
Bromomethane	20	U	20.4	102		24.1	121		17	*	15	70-130
Chloroethane	20	U	19.4	97		19.3	96		1		15	70-130
t-Butyl alcohol (TBA)	100	U	106.5	107		107.2	107		1		15	70-130
Trichlorofluoromethane	20	U	20.0	100		20.0	100		0		15	70-130
Diethyl ether	20	U	17.8	89		16.5	83		7		15	70-130
Acetone	100	13	105.6	93		100.8	88		5		15	70-130
1,1-Dichloroethene	20	U	19.8	99		19.6	98		1		15	70-130
Di-isopropyl ether (DIPE)	20	U	19.9	99		20.1	100		1		15	70-130
Methylene Chloride	20	U	18.4	92		18.5	93		1		15	70-130
Carbon Disulfide	20	U	17.2	86		17.4	87		1		15	70-130
Methyl-tert-butyl ether (MTBE)	40	U	41.0	102		41.4	104		1		15	70-130
trans-1,2-Dichloroethene	20	U	18.9	94		19.2	96		2		15	70-130
1,1-Dichloroethane	20	U	19.5	97		19.5	97		0		15	70-130
Methyl ethyl ketone	100	U	103.0	103		99.1	99		4		15	70-130
Ethyl t-butyl ether (ETBE)	20	U	19.7	99		20.1	100		2		15	70-130
2,2-Dichloropropane	20	U	21.4	107		21.4	107		0		15	70-130
cis-1,2-Dichloroethene	20	U	20.6	103		20.7	103		0		15	70-130
t-Amyl methyl ether (TAME)	20	U	20.3	101		20.7	103		2		15	70-130
Chloroform	20	U	20.3	101		20.6	103		1		15	70-130
Bromochloromethane	20	U	20.4	102		20.1	101		1		15	70-130
Tetrahydrofuran	20	U	19.5	97		19.7	98		1		15	70-130
1,1,1-Trichloroethane	20	U	20.6	103		20.9	104		2		15	70-130
1,1-Dichloropropene	20	U	19.9	99		20.4	102		2		15	70-130
Carbon Tetrachloride	20	U	21.6	108		22.2	111		3		15	70-130
1,2-Dichloroethane	20	U	19.5	97		19.5	97		0		15	70-130
Benzene	20	U	19.6	98		19.7	98		0		15	70-130
Trichloroethene	20	U	20.2	101		20.4	102		1		15	70-130
1,2-Dichloropropane	20	U	20.1	101		20.4	102		2		15	70-130
Bromodichloromethane	20	U	21.8	109		22.0	110		1		15	70-130
Dibromomethane	20	U	20.7	103		21.0	105		1		15	70-130
2-Hexanone	100	U	108.8	109		106.3	106		2		15	70-130
Methyl isobutyl ketone	100	U	106.4	106		103.2	103		3		15	70-130
cis-1,3-Dichloropropene	20	U	20.8	104		21.1	105		1		15	70-130
Toluene	20	U	19.9	99		20.0	100		1		15	70-130
trans-1,3-Dichloropropene	20	U	19.7	98		19.9	100		1		15	70-130
1,1,2-Trichloroethane	20	U	20.8	104		20.6	103		1		15	70-130
1,3-Dichloropropane	20	U	20.9	105		21.1	105		1		15	70-130
Tetrachloroethene	20	U	20.1	100		20.6	103		3		15	70-130
Dibromochloromethane	20	U	22.4	112		22.7	113		1		15	70-130
1,2-Dibromoethane	20	U	20.8	104		20.9	104		1		15	70-130
Chlorobenzene	20	U	19.6	98		19.9	100		1		15	70-130
1,1,1,2-Tetrachloroethane	20	U	21.4	107		22.1	111		3		15	70-130
Ethylbenzene	20	U	19.5	98		19.8	99		2		15	70-130
m,p-Xylene	40	U	38.7	97		39.3	98		2		15	70-130

analytics environmental laboratory LLC

VOLATILE ORGANIC
MATRIX SPIKE/DUPLICATE
PERCENT RECOVERY

Project #: 218736.002 Quote #: SDG#: 56637

Site ID: GP Old Town Mill Location: ME

Instrument ID: B
GC Column: RTX-502.2 ID: 0.25 (mm)

Heated Purge: (Y/N) N

Matrix Spike/Duplicate Sample ID: 56637-2

COMPOUND	SPIKE ADDED (ug/L)	SAMPLE AMT (ug/L)	MS CONC. (ug/L)	MS REC %	#	MSD CONC. (ug/L)	MSD REC %	#	RPD	#	QC LIMITS RPD	QC LIMITS % REC.
o-Xylene	20	U	20.1	100		20.5	102		2		15	70-130
Styrene	20	U	19.1	96		19.4	97		1		15	70-130
Bromoform	20	U	21.6	108		22.1	110		2		15	70-130
Isopropylbenzene	20	U	18.4	92		18.7	94		2		15	70-130
1,1,2,2-Tetrachloroethane	20	U	20.3	101		19.7	98		3		15	70-130
1,2,3-Trichloropropane	20	U	20.3	102		19.9	100		2		15	70-130
n-Propylbenzene	20	U	19.0	95		19.2	96		1		15	70-130
Bromobenzene	20	U	19.4	97		19.7	98		1		15	70-130
1,3,5-Trimethylbenzene	20	U	18.9	94		19.2	96		2		15	70-130
2-Chlorotoluene	20	U	19.7	99		19.8	99		0		15	70-130
4-Chlorotoluene	20	U	18.9	94		19.2	96		2		15	70-130
tert-butylbenzene	20	U	20.0	100		20.4	102		2		15	70-130
1,2,4-Trimethylbenzene	20	U	18.8	94		19.2	96		2		15	70-130
sec-butylbenzene	20	U	19.7	98		20.2	101		3		15	70-130
p-isopropyltoluene	20	2	20.9	95		21.5	98		3		15	70-130
1,3-Dichlorobenzene	20	U	18.8	94		19.1	95		2		15	70-130
1,4-Dichlorobenzene	20	U	18.6	93		19.1	95		2		15	70-130
n-butylbenzene	20	U	18.8	94		19.5	98		4		15	70-130
1,2-Dichlorobenzene	20	U	19.2	96		19.6	98		2		15	70-130
1,2-Dibromo-3-chloropropane	20	U	22.5	112		21.4	107		5		15	70-130
1,2,4-Trichlorobenzene	20	U	18.2	91		19.1	95		5		15	70-130
Hexachlorobutadiene	20	U	20.1	101		21.9	109		8		15	70-130
Naphthalene	20	U	19.8	99		19.9	99		0		15	70-130
1,2,3-Trichlorobenzene	20	U	18.2	91		19.3	96		5		15	70-130

Column to be used to flag recovery and RPD values with an asterisk

* Values outside of QC limits
RPD: 1 out of 71 outside limits
Spike Recovery: 0 out of 142 outside limits

Comments: _____

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VOLATILE ORGANIC

LABORATORY CONTROL SAMPLE/DUPLICATE

PERCENT RECOVERY

Project #: 218736.002 Quote #: SDG#: 56637

Site ID: GP Old Town Mill Location: ME

Instrument ID: B

GC Column: RTX-502.2 ID: 0.25 (mm)

Heated Purge: (Y/N) N

Laboratory Control Sample/Duplicate Sample ID: LS05306C/LS05306C2

COMPOUND	SPIKE ADDED (ug/kg)	SAMPLE AMT (ug/kg)	MS CONC. (ug/kg)	MS REC %	#	MSD CONC. (ug/kg)	MSD REC %	#	RPD	#	QC LIMITS RPD	QC LIMITS % REC.
Dichlorodifluoromethane	2000	U	1209	60		1135	57		6		25	49-82
Chloromethane	2000	U	1859	93		1731	87		7		25	75-125
Vinyl Chloride	2000	U	1414	71	*	1332	67	*	6		25	75-125
Bromomethane	2000	U	1951	98		2137	107		9		25	75-125
Chloroethane	2000	U	1637	82		1410	71	*	15		25	75-125
t-Butyl alcohol (TBA)	10000	U	7957	80		9103	91		13		25	60-140
Trichlorofluoromethane	2000	U	1807	90		1635	82		10		25	75-125
Diethyl ether	2000	U	1616	81		1505	75		7		25	75-125
Acetone	5000	U	4636	93		4322	86		7		25	75-125
1,1-Dichloroethene	2000	U	1604	80		1406	70	*	13		25	75-125
Di-isopropyl ether (DIPE)	2000	U	1737	87		1580	79		9		25	75-125
Methylene Chloride	2000	U	1726	86		1621	81		6		25	75-125
Carbon Disulfide	2000	U	1645	82		1518	76		8		25	75-125
Methyl-tert-butyl ether (MTBE)	2000	U	1706	85		1624	81		5		25	75-125
trans-1,2-Dichloroethene	2000	U	1752	88		1597	80		9		25	75-125
1,1-Dichloroethane	2000	U	1825	91		1632	82		11		25	75-125
Methyl ethyl ketone	5000	U	4306	86		4606	92		7		25	60-140
Ethyl t-butyl ether (ETBE)	2000	U	1778	89		1642	82		8		25	75-125
2,2-Dichloropropane	2000	U	2408	120		1435	72	*	51	*	25	75-125
cis-1,2-Dichloroethene	2000	U	1584	79		1711	86		8		25	75-125
t-Amyl methyl ether (TAME)	2000	U	1750	88		1664	83		5		25	75-125
Chloroform	2000	U	2252	113		1564	78		36	*	25	75-125
Bromochloromethane	2000	U	1662	83		1647	82		1		25	75-125
Tetrahydrofuran	2000	U	1708	85		1899	95		11		25	60-140
1,1,1-Trichloroethane	2000	U	1829	91		1639	82		11		25	75-125
1,1-Dichloropropene	2000	U	1803	90		1618	81		11		25	75-125
Carbon Tetrachloride	2000	U	1933	97		1725	86		11		25	75-125
1,2-Dichloroethane	2000	U	1779	89		1643	82		8		25	75-125
Benzene	2000	U	1666	83		1517	76		9		25	75-125
Trichloroethene	2000	U	1776	89		1733	87		2		25	75-125
1,2-Dichloropropane	2000	U	1792	90		1686	84		6		25	75-125
Bromodichloromethane	2000	U	1908	95		1801	90		6		25	75-125
Dibromomethane	2000	U	1731	87		1677	84		3		25	75-125
2-Hexanone	5000	U	4318	86		4437	89		3		25	75-125
Methyl isobutyl ketone	5000	U	4425	88		4503	90		2		25	75-125
cis-1,3-Dichloropropene	2000	U	1831	92		1768	88		3		25	75-125
Toluene	2000	U	1793	90		1711	86		5		25	75-125
trans-1,3-Dichloropropene	2000	U	1824	91		1758	88		4		25	75-125
1,1,2-Trichloroethane	2000	U	1786	89		1721	86		4		25	75-125
1,3-Dichloropropane	2000	U	1777	89		1761	88		1		25	75-125
Tetrachloroethene	2000	U	1880	94		1794	90		5		25	75-125
Dibromochloromethane	2000	U	2008	100		1933	97		4		25	75-125
1,2-Dibromoethane	2000	U	1769	88		1819	91		3		25	75-125
Chlorobenzene	2000	U	1792	90		1734	87		3		25	75-125
1,1,1,2-Tetrachloroethane	2000	U	1851	93		1707	85		8		25	75-125
Ethylbenzene	2000	U	1761	88		1665	83		6		25	75-125
m,p-Xylene	4000	U	3667	92		3449	86		6		25	75-125

analytics environmental laboratory LLC

VOLATILE ORGANIC

LABORATORY CONTROL SAMPLE/DUPLICATE

PERCENT RECOVERY

Project #: 218736.002 Quote #: SDG#: 56637

Site ID: GP Old Town Mill Location: ME

Instrument ID: B

GC Column: RTX-502.2 ID: 0.25 (mm)

Heated Purge: (Y/N) N

Laboratory Control Sample/Duplicate Sample ID: LS05306C/LS05306C2

COMPOUND	SPIKE ADDED (ug/kg)	SAMPLE AMT (ug/kg)	MS CONC. (ug/kg)	MS REC %	#	MSD CONC. (ug/kg)	MSD REC %	#	RPD	#	QC LIMITS RPD	QC LIMITS % REC.
o-Xylene	2000	U	1828	91		1764	88		4		25	75-125
Styrene	2000	U	1797	90		1719	86		4		25	75-125
Bromoform	2000	U	1934	97		1908	95		1		25	75-125
Isopropylbenzene	2000	U	1800	90		1703	85		6		25	75-125
1,1,2,2-Tetrachloroethane	2000	U	1703	85		1767	88		4		25	75-125
1,2,3-Trichloropropane	2000	U	1705	85		1758	88		3		25	75-125
n-Propylbenzene	2000	U	1895	95		1843	92		3		25	75-125
Bromobenzene	2000	U	1884	94		1823	91		3		25	75-125
1,3,5-Trimethylbenzene	2000	U	1864	93		1775	89		5		25	75-125
2-Chlorotoluene	2000	U	1831	92		1641	82		11		25	75-125
4-Chlorotoluene	2000	U	1813	91		1782	89		2		25	75-125
tert-butylbenzene	2000	U	1948	97		1749	87		11		25	75-125
1,2,4-Trimethylbenzene	2000	U	1828	91		1734	87		5		25	75-125
sec-butylbenzene	2000	U	1848	92		1742	87		6		25	75-125
p-isopropyltoluene	2000	U	1867	93		1817	91		3		25	75-125
1,3-Dichlorobenzene	2000	U	1887	94		1832	92		3		25	75-125
1,4-Dichlorobenzene	2000	U	1807	90		1764	88		2		25	75-125
n-butylbenzene	2000	U	1726	86		1693	85		2		25	75-125
1,2-Dichlorobenzene	2000	U	1757	88		1746	87		1		25	75-125
1,2-Dibromo-3-chloropropane	2000	U	1553	78		1682	84		8		25	75-125
1,2,4-Trichlorobenzene	2000	U	1792	90		1809	90		1		25	75-125
Hexachlorobutadiene	2000	U	1930	97		1818	91		6		25	75-125
Naphthalene	2000	U	1665	83		1851	93		11		25	75-125
1,2,3-Trichlorobenzene	2000	U	1825	91		1852	93		1		25	75-125

Column to be used to flag recovery and RPD values with an asterisk

* Values outside of QC limits

RPD: 2 out of 71 outside limits

Spike Recovery: 5 out of 142 outside limits

Comments: _____

_____

SS only V6010 Exp: 5/29/06
IS only V6011 Exp:
SS+IS V6012 Exp:

GC/MS Volatiles Instrument: C

052206-C

Voltage 1906
Threshold 100

Tube	Method	Sample	Vol	DF	Comments	File #	LB	pH
		W-C-052206-1-MT						
1	V805226C	BFB 50ng 2ml V6003	5mL		✓	C59021T	MT	NA
2		0.2 ppb .05ml V6017 V6007			✓	22si		
3		1 ppb .25ml			✓	23si		
4		7 ppb .50ml			✓	24si		
5		5 ppb 1.25ml			✓	25si		
6		20 ppb 5ml			✓	26si		
7		50 ppb 12.5ml			✓	27si		
8		100 ppb 25ml			✓	28si		
9		Blank				29		
10		L805226B 5ml V6018 V6014			RR - Cont. cell	30		
11		Blank			Cont cell? / CO issues	31		
12		L805226B2 5ml V6018 V6014			✓ Acrolein ↓	32		
13		L805226B3			✓ Acrolein still ↓	33		
14		Bromoform 5ml V6004			✓	34		
15		B805226B			✓	35		
16		56560-6			✓	36		<2
1		56475-7			✓	37		<2
2		~~56560-1~~ 56475-7, ~~500ml~~ 1mL			✓	38		<2
3	56560-1	-2	5mL		✓	39		<2
4	-2	-3			✓	40		<2
5	-3	-4			✓	41		<2
6	-4	-5			✓	42		<2
7		Blank			Cont cell	43		NA
8		Blank			Cont cell	44		
9		56560-5			Out of 12 hr window - RR 5mL	45		<2

MT 5/22/06

MT 5/23/06

Revision: 04/30/02

GC/MS Volatiles Instrument: B

IS only — Exp: —

SS+IS V6020 Exp: 6/7/06

052606-B

Voltage 2071

Threshold 100

Tube	Method	Sample	Vol	DF	Comments	File #	LB	pH
		w-B-052606-1-MT						
1	V8052606B	BFB 50ng 2ml V6023	5mL	✓		B3640TT	MT	NA
2		1 ppb .25ml V6017 V6022		✓		08si		
3		2 ppb .50ml		✓		09si		
4		5 ppb 1.25ml		✓		10si		
5		20 ppb 5ml		✓		11si		
6		50 ppb 12.5ml		✓		12si		
7		100 ppb 25ml		✓		13si		
8		Blank				14		
9		L8052606B 5ml V6018 V6014		✓		15Q		
10		L8052606B2		✓		16Q		
11		Bromoform 5ml V6024		✓		17		
12		B8052606B		✓		18B		
13		56594-4		✓		19		
14		-1		✓		20		
15		-2		✓		21		
16		-3		✓		22		
17		56603-1		✓		23		
18		-1MS		✓		24sp		
19		-1MSD		✓		25sp		
20		-2		✓		26		
21		-3		✓		27		
22		-4		✓		28		
23		MT 5/26/06 56602-1		✓		29		
24		-6 -2		✓		30		
25		BFB 50ng 2ml V6023		✓		31		NA
26		25 ppb 5ml V6017 V6022		✓		32		
27		L8052606B3 5ml V6018 V6014		✓		33		
28		L8052606B4		✓		34		
29		Blank				35		
30		B8052606B2		✓		36		

cont on next page.

Revision: 04/30/02

GC/MS Volatiles Instrument: B

IS only Exp:

SS+IS V6020 Exp: 6-7-06

Voltage 2071

Threshold 100

Jm 053001-B

Tube	Method	Sample	Vol	DF	Comments	File #	LB	pH
	W-B-053001-1-JR							
2	V805266B	BFB 50ng 2mL V6023	5mL	✓		B36458T	Jm	
3		20PPB 5mL V6017 V6022		✓		59	SC	
4		L805306B V6014 V6018		✓		60	Q	
5		Bromoform		✓		61		
6		B805306B		✓		62	B	
7		56637-29	5mL	✓		63		<2
8		56638-5		✓		64		<2
9		↓ -9		✓		65		<2
10		56603-25		✓		66		<2
11		56618-5	1mL	✓		67		<2
12		↓ -9	500mL	✓		68		<2
13		56637-1	5mL	✓		69		<2
14		-2		✓		70		<2
15		-2,MS		✓	} BM RPD 17%	71	SP	<2
16		-2,MSD		✓		72	SP	<2
17		-3		✓		73		<2
18		-5		✓		74		Unpres.
19		-20		✓		75		<2
20		-21		✓		76		<2
21		-24		✓		77		Unpres.
22		-25		✓		78		
23		-26		✓		79		
24		-27		✓		80		
25		-28		✓		81		
26		BFB 50ng 2mL V6023	5mL	✓		82	T	
27		20PPB 5mL V6017 V6022		✓		83	SC	
28		L805306B2 V6014 V6018		✓		84	Q	
29		Blank				85		
30		B805306B2		✓		86	B	
31		56629-5	5mL	✓		87		<2

☒ 053001B1.XLS

☐ 052601B1.XLS

Revision: 04/30/02

GC/MS Volatiles Instrument: C

IS only V6011 Exp: 5/29/04
SS+IS V6012 Exp: 5/29/04
Voltage 1906
Threshold 100

053006-C

ube	Method	Sample	Vol	DF	Comments		File #	LB	pH
		W-C-053006-1-SC							
2	V505226C	BFB 50ng 2ul V6023	5mL		✓		C59117T	SL	N/A
3		20ppb 5ul ✓V6017 ✓V6022	↓		✓		1185C		
4		LS05306C ✓V6018 ✓V6024	50ul/5mL		✓		119Q		
5		LS05306CZ ↓	↓		✓		120Q		
6		Bromofluorm 5uL V6024	5mL		✓		121		
7		MB05306C	50ul/5mL		✓		122B		
15		56637-4			✓	4.39	123		
16		-14			✓ DSS, Re-inject	4.10	124		
6		-15			✓	3.08	125		
7		-16			✓	4.98	126		
8		-17			*RR WL C10	4.78	127		
9		-19			✓	4.37	128		
10		-9			✓*	5.00	129		
11		-10			✓*	5.00	130		
12		-11			✓*	5.00	131		
13		-18			✓ DSS, Re-inject	5.60	132		
14		-18			10mL F.V. DSS, LK	5.00	133		
15		-14, RR			DSS, Confirms	4.10	134		
16		-17, RR			✓	4.78	135		
1		-18, RR			DSS, Confirms	5.60	136		
2		-18, RR			✓ 10mL F.V.	5.00	137		
3		56576-1			25mL F.V. DSS, RR	01.01	138		
4	↓	56599-1	↓		DSS, out of window	9.37	139	↓	↓

5/31/06 SL

* these samples had no weight recorded, Assuming 5.00g

Revision: 04/30/02

analytics environmental laboratory LLC

VOLATILE ORGAN
INSTRUMENT PERFORMANCE CHECK
BROMOFLUOROBENZENE (BFB)

Project #: 218736.002 Quote #: SDG#: 56637

BFB File ID: C59021T.D Site ID: GP Old Town Mill Location: ME

Instrument ID: C BFB Injection Date: 5/22/06

GC Column: RTX-502.2 ID: 0.25(mm) BFB Injection Time: 1030

Heated Purge: (Y/N) N

m/e	ION ABUNDANCE CRITERIA	%RELATIVE ABUNDANCE
50	15.0 - 40.0% of mass 95	21.2
75	30.0 - 60.0% of mass 95	45.8
95	Base peak, 100% relative abundance	100.0
96	5.0 - 9.0% of mass 95	6.5
173	Less than 2.0% of mass 174	0.0 (0.0)1
174	Greater than 50% of m	99.0
175	5.0 - 9.0% of mass 174	7.0 (7.1)1
176	>95.0 - <101.0% of mass 174	94.7 (95.7)1
177	5.0 - 9.0% of mass 176	6.2 (6.5)2

1-Value is % mass 174 2-Value is % mass 176

This check applies to the following SAMPLES, MS, MSD, BLANKS and STANDARDS:

	SAMPLE NO.	LAB SAMPLE	LAB FILE ID	DATE ANALYZED	TIME ANALYZED
01	0.2 PPB	0.2 PPB	C59022SI.D	05/22/06	10:30
02	1 PPB	1 PPB	C59023SI.D	05/22/06	11:24
03	2 PPB	2 PPB	C59024SI.D	05/22/06	11:57
04	5 PPB	5 PPB	C59025SI.D	05/22/06	12:30
05	20 PPB	20 PPB	C59026SI.D	05/22/06	13:07
06	50 PPB	50 PPB	C59027SI.D	05/22/06	13:07
07	100 PPB	100 PPB	C59028SI.D	05/22/06	13:50
08	ZZZZZ	ZZZZZ	C59029.D	05/22/06	14:38
09	ZZZZZ	ZZZZZ	C59030Q.D	05/22/06	15:13
10	ZZZZZ	ZZZZZ	C59031.D	05/22/06	15:46
11	ZZZZZ	ZZZZZ	C59032Q.D	05/22/06	16:15
12	ZZZZZ	ZZZZZ	C59033Q.D	05/22/06	16:48
13	ZZZZZ	ZZZZZ	C59034.D	05/22/06	16:48
14	ZZZZZ	ZZZZZ	C59035B.D	05/22/06	17:35
15	ZZZZZ	ZZZZZ	C59036.D	05/22/06	18:14
16	ZZZZZ	ZZZZZ	C59037.D	05/22/06	18:47
17	ZZZZZ	ZZZZZ	C59038.D	05/22/06	18:52
18	ZZZZZ	ZZZZZ	C59039.D	05/22/06	19:26
19	ZZZZZ	ZZZZZ	C59040.D	05/22/06	19:59
20	ZZZZZ	ZZZZZ	C59041.D	05/22/06	20:32
21	ZZZZZ	ZZZZZ	C59042.D	05/22/06	21:05
22	ZZZZZ	ZZZZZ	C59043.D	05/22/06	21:43
23	ZZZZZ	ZZZZZ	C59044.D	05/22/06	22:16
24	ZZZZZ	ZZZZZ	C59045.D	05/22/06	22:50

analytics environmental laboratory LLC

VOLATILE ORGAN
INSTRUMENT PERFORMANCE CHECK
BROMOFLUOROBENZENE (BFB)

Project #:	218736.005	Quote #:		SDG#:	56637
BFB File ID:	B36407T.D	Site ID:	GP Old Town Mill	Location:	ME
Instrument ID:	B			BFB Injection Date:	5/26/06
GC Column:	RTX-502.2	ID: 0.25(mm)		BFB Injection Time:	0956
				Heated Purge: (Y/N)	N

m/e	ION ABUNDANCE CRITERIA	%RELATIVE ABUNDANCE	
50	15.0 - 40.0% of mass 95	26.2	
75	30.0 - 60.0% of mass 95	54.3	
95	Base peak, 100% relative abundance	100.0	
96	5.0 - 9.0% of mass 95	6.8	
173	Less than 2.0% of mass 174	0.0 (	0.0)1
174	Greater than 50% of m	74.2	
175	5.0 - 9.0% of mass 174	6.2 (	8.3)1
176	>95.0 - <101.0% of mass 174	72.0 (	97.0)1
177	5.0 - 9.0% of mass 176	4.6 (	6.4)2

1-Value is % mass 174 2-Value is % mass 176

This check applies to the following SAMPLES, MS, MSD, BLANKS and STANDARDS:

	SAMPLE NO.	LAB SAMPLE	LAB FILE ID	DATE ANALYZED	TIME ANALYZED
01	1 PPB	1 PPB	B36408SI.D	05/26/06	10:18
02	2 PPB	2 PPB	B36409SI.D	05/26/06	10:53
03	5 PPB	5 PPB	B36410SI.D	05/26/06	11:22
04	20 PPB	20 PPB	B36411SI.D	05/26/06	11:52
05	50 PPB	50 PPB	B36412SI.D	05/26/06	12:21
06	100 PPB	100 PPB	B36413SI.D	05/26/06	12:50
07	ZZZZZ	ZZZZZ	B36414.D	05/26/06	13:24
08	ZZZZZ	ZZZZZ	B36415Q.D	05/26/06	13:44
09	ZZZZZ	ZZZZZ	B36416Q.D	05/26/06	14:17
10	ZZZZZ	ZZZZZ	B36417.D	05/26/06	14:47
11	ZZZZZ	ZZZZZ	B36418B.D	05/26/06	15:16
12	ZZZZZ	ZZZZZ	B36419.D	05/26/06	15:46
13	ZZZZZ	ZZZZZ	B36420.D	05/26/06	16:15
14	ZZZZZ	ZZZZZ	B36421.D	05/26/06	16:44
15	ZZZZZ	ZZZZZ	B36422.D	05/26/06	17:13
16	ZZZZZ	ZZZZZ	B36423.D	05/26/06	17:43
17	ZZZZZ	ZZZZZ	B36424SP.D	05/26/06	18:12
18	ZZZZZ	ZZZZZ	B36425SP.D	05/26/06	18:41
19	ZZZZZ	ZZZZZ	B36426.D	05/26/06	19:11
20	ZZZZZ	ZZZZZ	B36427.D	05/26/06	19:40
21	ZZZZZ	ZZZZZ	B36428.D	05/26/06	20:09
22	ZZZZZ	ZZZZZ	B36429.D	05/26/06	20:38
23	ZZZZZ	ZZZZZ	B36430.D	05/26/06	21:08
24					

analytics environmental laboratory LLC

VOLATILE ORGAN
INSTRUMENT PERFORMANCE CHECK
BROMOFLUOROBENZENE (BFB)

Project #: 218736.002 Quote #: SDG#: 56637

BFB File ID: C59117T.D Site ID: GP Old Town Mill Location: ME

Instrument ID: C BFB Injection Date: 5/30/06

GC Column: RTX-502.2 ID: 0.25(mm) BFB Injection Time: 0818

Heated Purge: (Y/N) N

m/e	ION ABUNDANCE CRITERIA	%RELATIVE ABUNDANCE
50	15.0 - 40.0% of mass 95	22.6
75	30.0 - 60.0% of mass 95	46.9
95	Base peak, 100% relative abundance	100.0
96	5.0 - 9.0% of mass 95	7.3
173	Less than 2.0% of mass 174	0.0 (0.4)1
174	Greater than 50% of m	98.0
175	5.0 - 9.0% of mass 174	4.5 (7.4)1
176	>95.0 - <101.0% of mass 174	57.6 (97.6)1
177	5.0 - 9.0% of mass 176	3.8 (6.7)2

1-Value is % mass 174 2-Value is % mass 176

This check applies to the following SAMPLES, MS, MSD, BLANKS and STANDARDS:

	SAMPLE NO.	LAB SAMPLE	LAB FILE ID	DATE ANALYZED	TIME ANALYZED
01	20 PPB	20 PPB	C59118SC.D	05/30/06	8:45
02	LS05306C	LS05306C	C59119Q.D	05/30/06	9:10
03	LS05306C2	LS05306C2	C59120Q.D	05/30/06	9:43
04	ZZZZZ	ZZZZZ	C59121.D	05/30/06	10:16
05	MB05306C	MB05306C	C59122B.D	05/30/06	10:55
06	56637-4	56637-4	C59123.D	05/30/06	11:30
07	56637-14	56637-14	C59124.D	05/30/06	12:03
08	56637-15	56637-15	C59125.D	05/30/06	12:36
09	56637-16	56637-16	C59126.D	05/30/06	13:14
10	ZZZZZ	ZZZZZ	C59127.D	05/30/06	13:26
11	56637-19	56637-19	C59128.D	05/30/06	14:17
12	56637-9	56637-9	C59129.D	05/30/06	14:46
13	56637-10	56637-10	C59130.D	05/30/06	15:21
14	56637-11	56637-11	C59131.D	05/30/06	15:55
15	ZZZZZ	ZZZZZ	C59132.D	05/30/06	15:55
16	ZZZZZ	ZZZZZ	C59133.D	05/30/06	17:38
17	56637-14, RR	56637-14, RR	C59134.D	05/30/06	17:59
18	56637-17, RR	56637-17, RR	C59135.D	05/30/06	18:33
19	ZZZZZ	ZZZZZ	C59136.D	05/30/06	19:06
20	56637-18, RR,, 10ML FV	56637-18, RR,, 10ML FV	C59137.D	05/30/06	19:10
21	ZZZZZ	ZZZZZ	C59138.D	05/30/06	20:05
22	ZZZZZ	ZZZZZ	C59139.D	05/30/06	20:42
23					
24					

VOLATILE ORGAN
INSTRUMENT PERFORMANCE CHECK
BROMOFLUOROBENZENE (BFB)

Project #:	218736.002	Quote #:		SDG#:	56637
BFB File ID:	B36458T.D	Site ID:	GP Old Town Mill	Location:	ME
Instrument ID:	B			BFB Injection Date:	5/30/06
GC Column:	RTX-502.2	ID: 0.25(mm)		BFB Injection Time:	0858
				Heated Purge: (Y/N)	N

m/e	ION ABUNDANCE CRITERIA	%RELATIVE ABUNDANCE
50	15.0 - 40.0% of mass 95	25.7
75	30.0 - 60.0% of mass 95	52.4
95	Base peak, 100% relative abundance	100.0
96	5.0 - 9.0% of mass 95	6.8
173	Less than 2.0% of mass 174	0.0 (0.0)1
174	Greater than 50% of m	69.1
175	5.0 - 9.0% of mass 174	5.7 (8.2)1
176	>95.0 - <101.0% of mass 174	68.9 (99.8)1
177	5.0 - 9.0% of mass 176	4.7 (6.8)2

1-Value is % mass 174 2-Value is % mass 176

This check applies to the following SAMPLES, MS, MSD, BLANKS and STANDARDS:

	SAMPLE NO.	LAB SAMPLE	LAB FILE ID	DATE ANALYZED	TIME ANALYZED
01	20 PPB	20 PPB	B36459SC.D	05/30/06	9:35
02	L805306B	L805306B	B36460Q.D	05/30/06	9:52
03	ZZZZZ	ZZZZZ	B36461.D	05/30/06	10:22
04	B805306B	B805306B	B36462B.D	05/30/06	10:51
05	56637-29	56637-29	B36463.D	05/30/06	11:20
06	ZZZZZ	ZZZZZ	B36464.D	05/30/06	11:50
07	ZZZZZ	ZZZZZ	B36465.D	05/30/06	12:19
08	ZZZZZ	ZZZZZ	B36466.D	05/30/06	12:57
09	ZZZZZ	ZZZZZ	B36467.D	05/30/06	13:26
10	ZZZZZ	ZZZZZ	B36468.D	05/30/06	13:56
11	56637-1	56637-1	B36469.D	05/30/06	14:25
12	56637-2	56637-2	B36470.D	05/30/06	14:54
13	56637-2,MS	56637-2,MS	B36471SP.D	05/30/06	15:24
14	56637-2,MSD	56637-2,MSD	B36472SP.D	05/30/06	15:53
15	56637-3	56637-3	B36473.D	05/30/06	16:22
16	56637-5	56637-5	B36474.D	05/30/06	16:51
17	56637-20	56637-20	B36475.D	05/30/06	17:21
18	56637-21	56637-21	B36476.D	05/30/06	17:50
19	56637-24	56637-24	B36477.D	05/30/06	18:19
20	56637-25	56637-25	B36478.D	05/30/06	18:48
21	56637-26	56637-26	B36479.D	05/30/06	19:18
22	56637-27	56637-27	B36480.D	05/30/06	19:47
23	56637-28	56637-28	B36481.D	05/30/06	20:16
24					

Acq. Method File	**V805226C**
Level 1 File Name	C59023SI.D
Level 2 File Name	C59024SI.D
Level 3 File Name	C59025SI.D
Level 4 File Name	C59026SI.D
Level 5 File Name	C59027SI.D
Level 6 File Name	C59028SI.D
Level 7 File Name	C59022SI.D

5-23-06

Name	Level 1 RF	Level 2 RF	Level 3 RF	Level 4 RF	Level 5 RF	Level 6 RF	Level 7 RF	AVG RF	RF Rel. STDEV
Dichlorodifluoromethane	0.707	0.532	0.579	0.581	0.592	0.618	-1.000	0.373	9.74%
Chloromethane	0.651	0.531	0.571	0.613	0.678	0.770	-1.000	0.402	13.30%
Vinyl Chloride	0.779	0.600	0.649	0.619	0.585	0.622	-1.000	0.408	10.95%
Bromomethane	0.323	0.238	0.366	0.398	0.425	-1.000	0.000	0.107	20.93%
Chloroethane	0.459	0.319	0.324	0.310	0.313	0.360	-1.000	0.155	16.52%
t-Butyl alcohol (TBA)	0.069	0.060	0.054	0.052	0.057	-1.000	0.000	-0.101	11.13%
Trichlorofluoromethane	1.489	1.043	1.163	1.154	1.172	1.256	-1.000	0.897	12.49%
Diethyl ether	0.212	0.241	0.206	0.196	0.190	0.200	-1.000	0.035	8.74%
1,1,2-Trichlorotrifluoroethane	0.612	0.572	0.510	0.515	0.511	0.538	-1.000	0.323	7.63%
Acrolein	0.083	0.065	0.075	0.077	0.051	-1.000	0.000	-0.093	18.15% NT
Acetone	0.195	0.188	0.170	0.176	0.183	-1.000	0.000	-0.012	5.26%
1,1-Dichloroethene	0.529	0.424	0.390	0.384	0.376	0.396	-1.000	0.214	13.84%
Methyl iodide	0.700	0.667	0.655	0.835	0.713	0.812	-1.000	0.483	10.34%
Di-isopropyl ether (DIPE)	2.063	1.879	1.705	1.739	1.726	1.832	-1.000	1.421	7.42%
Methylene Chloride	0.525	0.427	0.403	0.391	0.410	-1.000	0.000	0.165	12.53%
Carbon Disulfide	1.326	1.364	1.250	1.320	1.342	1.424	-1.000	1.004	4.27%
Acrylonitrile	0.188	0.197	0.180	0.178	0.195	-1.000	0.000	-0.009	4.54%
Methyl-tert-butyl ether (MTBE)	1.728	1.616	1.417	1.408	1.403	1.503	-1.000	1.154	8.83%
trans-1,2-Dichloroethene	1.067	0.982	0.898	0.870	0.900	0.948	-1.000	0.666	7.64%
1,1-Dichloroethane	1.117	1.131	1.022	1.033	1.049	1.134	-1.000	0.784	4.79%
Vinyl acetate	1.403	1.258	0.990	1.108	1.181	1.163	-1.000	0.872	11.80%
Methyl ethyl ketone	0.067	0.042	0.049	0.046	0.045	0.050	-1.000	-0.100	17.26%
Ethyl t-butyl ether (ETBE)	2.133	1.935	1.771	1.810	1.834	1.991	-1.000	1.497	7.10%
2,2-Dichloropropane	1.166	0.834	0.829	0.772	0.739	1.216	-1.000	0.651	22.56%
cis-1,2-Dichloroethene	0.621	0.643	0.608	0.618	0.600	0.583	-1.000	0.382	3.30%
t-Amyl methyl ether (TAME)	1.529	1.379	1.238	1.268	1.253	1.301	-1.000	0.995	8.32%
DB Holder #28	0.991	0.721	0.726	0.738	0.730	0.786	-1.000	0.527	13.42%
Chloroform	1.156	1.056	0.959	0.991	0.988	1.414	-1.000	0.795	15.70% 10K CCC
Bromochloromethane	0.368	0.360	0.317	0.328	0.316	0.326	-1.000	0.145	6.66%
Tetrahydrofuran	0.056	0.059	0.055	0.053	0.057	-1.000	0.000	-0.103	4.12%
1,1,1-Trichloroethane	1.025	0.916	0.868	0.899	0.905	0.979	-1.000	0.656	6.29%
1,1-Dichloropropene	0.873	0.797	0.716	0.734	0.725	0.755	-1.000	0.514	7.77%
Carbon Tetrachloride	0.771	0.743	0.718	0.786	0.801	0.886	-1.000	0.529	7.43%
d4-1,2-Dichloroethane	0.991	0.721	0.726	0.738	0.730	0.786	-1.000	0.527	13.42%
1,2-Dichloroethane	1.098	1.001	0.914	0.949	0.947	1.045	-1.000	0.708	7.02%
Benzene	2.104	1.923	1.747	1.742	1.652	1.716	2.587	1.924	17.15%
Trichloroethene	0.426	0.397	0.361	0.361	0.362	0.358	-1.000	0.181	7.41%
1,2-Dichloropropane	0.375	0.386	0.342	0.335	0.330	0.324	-1.000	0.156	7.35%
Methylmethacrylate	0.235	0.220	0.248	0.215	0.200	0.210	-1.000	0.047	7.79%
Bromodichloromethane	0.459	0.434	0.413	0.427	0.441	0.464	-1.000	0.234	4.36%
Dibromomethane	0.291	0.272	0.266	0.270	0.259	0.267	-1.000	0.089	4.02%
1,4-Dioxane	0.004	0.004	0.004	0.003	0.004	0.003	-1.000	-0.140	11.20%
2-Chloroethylvinyl ether	0.093	0.086	0.075	0.114	0.104	0.170	-1.000	-0.051	99.70% NT
2-Hexanone	0.146	0.140	0.127	0.119	0.120	0.117	-1.000	-0.033	9.63%
Methyl isobutyl ketone	0.147	0.143	0.131	0.124	0.118	0.118	-1.000	-0.031	9.60%
cis-1,3-Dichloropropene	0.518	0.523	0.482	0.500	0.506	0.512	-1.000	0.292	2.89%
d8-Toluene	1.276	0.991	1.063	1.053	1.046	1.050	1.522	1.143	16.64%
Toluene	0.807	0.776	0.686	0.702	0.685	0.689	-1.000	0.478	7.37%
trans-1,3-Dichloropropene	0.522	0.493	0.459	0.460	0.464	0.493	-1.000	0.270	5.22%
1,1,2-Trichloroethane	0.257	0.265	0.232	0.231	0.229	0.228	-1.000	0.063	6.80%
1,3-Dichloropropane	0.516	0.495	0.452	0.442	0.435	0.441	-1.000	0.254	7.24%
Tetrachloroethene	0.475	0.467	0.413	0.417	0.412	0.428	-1.000	0.230	6.55%
Dibromochloromethane	0.400	0.361	0.333	0.372	0.392	0.409	-1.000	0.181	7.47%
1,2-Dibromoethane	0.442	0.426	0.389	0.392	0.384	0.404	-1.000	0.205	5.68%
Chlorobenzene	1.201	1.146	1.014	1.044	1.033	0.999	-1.000	0.777	7.59%
1,1,1,2-Tetrachloroethane	0.412	0.404	0.341	0.369	0.377	0.372	-1.000	0.182	6.75%
Ethylbenzene	1.811	1.639	1.465	1.503	1.504	1.403	-1.000	1.189	9.51%
m,p-Xylene	0.739	0.692	0.606	0.599	0.578	0.527	-1.000	0.392	12.52%
o-Xylene	0.673	0.675	0.566	0.573	0.581	0.549	-1.000	0.374	9.33%
Styrene	1.185	1.161	1.012	1.020	1.011	0.980	-1.000	0.767	8.28%

Bromoform	0.311	0.313	0.301	0.319	0.356	0.352	-1.000	0.136	7.09%
Isopropylbenzene	1.853	1.728	1.517	1.539	1.540	1.457	-1.000	1.233	9.46%
1,1,2,2-Tetrachloroethane	0.491	0.457	0.414	0.402	0.405	0.386	-1.000	0.222	9.39%
romofluorobenzene	0.651	0.487	0.471	0.461	0.468	0.452	-1.000	0.284	15.17%
1,2,3-Trichloropropane	0.349	0.322	0.315	0.291	0.298	0.282	-1.000	0.122	7.88%
trans-1,4-Dichloro-2-butene	0.143	0.135	0.125	0.119	0.121	0.114	-1.000	-0.035	8.59%
n-Propylbenzene	2.144	1.962	1.720	1.708	1.660	1.458	-1.000	1.379	13.63%
Bromobenzene	0.497	0.484	0.435	0.431	0.445	0.415	-1.000	0.244	7.16%
1,3,5-Trimethylbenzene	1.463	1.381	1.235	1.228	1.244	1.166	-1.000	0.960	8.69%
2-Chlorotoluene	1.405	1.234	1.142	1.145	1.136	1.084	-1.000	0.878	9.71%
4-Chlorotoluene	1.232	1.223	1.026	1.018	1.074	1.009	-1.000	0.798	9.44%
tert-butylbenzene	1.250	1.224	1.074	1.066	1.110	1.020	-1.000	0.820	8.22%
1,2,4-Trimethylbenzene	1.490	1.488	1.270	1.201	1.273	1.189	-1.000	0.988	10.35%
sec-butylbenzene	2.001	1.924	1.692	1.656	1.702	1.645	-1.000	1.374	8.61%
p-isopropyltoluene	1.539	1.540	1.332	1.325	1.349	1.289	-1.000	1.053	8.10%
1,3-Dichlorobenzene	0.855	0.804	0.716	0.714	0.730	0.703	-1.000	0.503	8.16%
1,4-Dichlorobenzene	1.703	1.663	1.579	1.552	1.556	1.498	-1.000	1.221	4.83%
n-butylbenzene	3.058	2.767	2.528	2.587	2.549	2.514	-1.000	2.143	7.97%
1,2-Dichlorobenzene	1.653	1.474	1.412	1.446	1.403	1.377	-1.000	1.109	6.86%
1,2-Dibromo-3-chloropropane	0.136	0.168	0.151	0.141	0.150	0.149	-1.000	-0.015	7.26%
1,2,4-Trichlorobenzene	1.393	1.200	1.145	1.179	1.194	1.216	-1.000	0.904	7.19%
Hexachlorobutadiene	0.679	0.595	0.512	0.581	0.609	0.621	-1.000	0.371	9.09%
Naphthalene	2.681	2.304	2.234	2.153	2.287	2.234	-1.000	1.842	8.07%
1,2,3-Trichlorobenzene	1.165	1.067	1.048	1.036	1.101	1.097	-1.000	0.788	4.29%
1,3,5-Trichlorobenzene	1.493	1.384	1.256	1.305	1.341	1.372	-1.000	1.022	5.95%
								AVG	**10.24%**

Response Factor Report Instr_C

Method : D:\HPCHEM\1\METHODS\V805226C.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Mon May 22 14:18:09 2006
Response via : Initial Calibration

NT 5/23/06

5.23.06

Calibration Files
1 =C59023SI.D 2 =C59024SI.D 5 =C59025SI.D
20 =C59026SI.D 50 =C59027SI.D 100 =C59028SI.D

		Compound	1	2	5	20	50	100	Avg	%RSD
1)	I	Pentafluorobenzene	----------	ISTD	----------					
2)	TM	Dichlorodifluoromet	0.707	0.532	0.579	0.581	0.592	0.618	0.601	9.74
3)	TMP	Chloromethane	0.651	0.531	0.571	0.613	0.678	0.770	0.636✓	13.30
4)	TMC	Vinyl Chloride	0.779	0.600	0.649	0.619	0.585	0.622	0.642	10.95✓
5)	TM	Bromomethane	0.323	0.238	0.366	0.398	0.425		0.350	20.93
6)	TM	Chloroethane	0.459	0.319	0.324	0.310	0.313	0.360	0.347	16.52
7)	TM	t-Butyl alcohol (TB		0.069	0.060	0.054	0.052	0.057	0.058#	11.13
8)	TM	Trichlorofluorometh	1.489	1.043	1.163	1.154	1.172	1.256	1.213	12.49
9)	TM	Diethyl ether	0.212	0.241	0.206	0.196	0.190	0.200	0.208	8.74
10)	T	1,1,2-Trichlorotrif	0.612	0.572	0.510	0.515	0.511	0.538	0.543	7.63
11)	T	Acrolein		0.083	0.065	0.075	0.077	0.051	0.070#	18.15 NT
12)	TM	Acetone		0.195	0.188	0.170	0.176	0.183	0.183	5.26
13)	TMC	1,1-Dichloroethene	0.529	0.424	0.390	0.384	0.376	0.396	0.416	13.84✓
14)	T	Methyl iodide	0.700	0.667	0.655	0.835	0.713	0.812	0.731	10.34
15)	TM	Di-isopropyl ether	2.063	1.879	1.705	1.739	1.726	1.832	1.824	7.42
16)	TM	Methylene Chloride		0.525	0.427	0.403	0.391	0.410	0.431	12.53
17)	TM	Carbon Disulfide	1.326	1.364	1.250	1.320	1.342	1.424	1.337	4.27
18)	T	Acrylonitrile		0.188	0.197	0.180	0.178	0.195	0.188	4.54
19)	TM	Methyl-tert-butyl e	1.728	1.616	1.417	1.408	1.403	1.503	1.513	8.83
20)	TM	trans-1,2-Dichloroe	1.067	0.982	0.898	0.870	0.900	0.948	0.944	7.64
21)	TMP	1,1-Dichloroethane	1.117	1.131	1.022	1.033	1.049	1.134	1.081✓	4.79
22)	T	Vinyl acetate	1.403	1.258	0.990	1.108	1.181	1.163	1.184	11.80
23)	TM	Methyl ethyl ketone	0.067	0.042	0.049	0.046	0.045	0.050	0.050#	17.26
24)	TM	Ethyl t-butyl ether	2.133	1.935	1.771	1.810	1.834	1.991	1.913	7.10
25)	TM	2,2-Dichloropropane	1.166	0.834	0.829	0.772	0.739	1.216	0.926	22.56
26)	TM	cis-1,2-Dichloroeth	0.621	0.643	0.608	0.618	0.600	0.583	0.612	3.30
27)	TM	t-Amyl methyl ether	1.529	1.379	1.238	1.268	1.253	1.301	1.328	8.32
28)	S	DB Holder #28	0.991	0.721	0.726	0.738	0.730	0.786	0.782	13.42
29)	TMC	Chloroform	1.156	1.056	0.959	0.991	0.988	1.414	1.094	15.70#✓
30)	TM	Bromochloromethane	0.368	0.360	0.317	0.328	0.316	0.326	0.336	6.66
31)	TM	Tetrahydrofuran		0.056	0.059	0.055	0.053	0.057	0.056#	4.12
32)	TM	1,1,1-Trichloroetha	1.025	0.916	0.868	0.899	0.905	0.979	0.932	6.29
33)	T	NO COMPOUND							0.000#	-1.00
34)	TM	1,1-Dichloropropene	0.873	0.797	0.716	0.734	0.725	0.755	0.767	7.77
35)	TM	Carbon Tetrachlorid	0.771	0.743	0.718	0.786	0.801	0.886	0.784	7.43
36)	S	d4-1,2-Dichloroetha	0.991	0.721	0.726	0.738	0.730	0.786	0.782	13.42
37)	TM	1,2-Dichloroethane	1.098	1.001	0.914	0.949	0.947	1.045	0.992	7.02
38)	TM	Benzene	2.104	1.923	1.747	1.742	1.652	1.716	1.924	17.15
39)	I	1,4-Difluorobenzene	----------	ISTD	----------					
40)	TM	Trichloroethene	0.426	0.397	0.361	0.361	0.362	0.358	0.377	7.41
41)	TMC	1,2-Dichloropropane	0.375	0.386	0.342	0.335	0.330	0.324	0.349	7.35✓
42)	T	Methylmethacrylate	0.235	0.220	0.248	0.215	0.200	0.210	0.221	7.79

(#) = Out of Range ### Number of calibration levels exceeded format ###

Response Factor Report Instr_C

Method : D:\HPCHEM\1\METHODS\V805226C.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Mon May 22 14:18:09 2006
Response via : Initial Calibration

Calibration Files
1 =C59023SI.D 2 =C59024SI.D 5 =C59025SI.D
20 =C59026SI.D 50 =C59027SI.D 100 =C59028SI.D

		Compound	1	2	5	20	50	100	Avg	%RSD
43)	T	Methyl acetate	0.319	0.292	0.323	0.265	0.244	0.256	0.283	11.77
44)	T	Cyclohexane	0.672	0.618	0.552	0.549	0.535	0.502	0.571	10.93
45)	TM	Bromodichloromethan	0.459	0.434	0.413	0.427	0.441	0.464	0.440	4.36
46)	TM	Dibromomethane	0.291	0.272	0.266	0.270	0.259	0.267	0.271	4.02
47)	T	1,4-Dioxane	0.004	0.004	0.004	0.003	0.004	0.003	0.004#	11.20
48)	T	Methylcyclohexane	0.556	0.511	0.461	0.460	0.455	0.453	0.483	8.71
49)	T	NO COMPOUND							0.000#	-1.00
50)	T	2-Chloroethylvinyl	0.093	0.086	0.075	0.114	0.104	0.170	0.107	31.66 Quad
51)	TM	2-Hexanone	0.146	0.140	0.127	0.119	0.120	0.117	0.128	9.63
52)	TM	Methyl isobutyl ket	0.147	0.143	0.131	0.124	0.118	0.118	0.130	9.60
53)	TM	cis-1,3-Dichloropro	0.518	0.523	0.482	0.500	0.506	0.512	0.507	2.89
54)	S	d8-Toluene	1.276	0.991	1.063	1.053	1.046	1.050	1.143	16.64
55)	TMC	Toluene	0.807	0.776	0.686	0.702	0.685	0.689	0.724	7.37✓
56)	TM	trans-1,3-Dichlorop	0.522	0.493	0.459	0.460	0.464	0.493	0.482	5.22
57)	T	NO COMPOUND							0.000#	-1.00
58)	TM	1,1,2-Trichloroetha	0.257	0.265	0.232	0.231	0.229	0.228	0.240	6.80
59)	TM	1,3-Dichloropropane	0.516	0.495	0.452	0.442	0.435	0.441	0.463	7.24
60)	TM	Tetrachloroethene	0.475	0.467	0.413	0.417	0.412	0.428	0.435	6.55
61)	TM	Dibromochloromethan	0.400	0.361	0.333	0.372	0.392	0.409	0.378	7.47
62)	TM	1,2-Dibromoethane	0.442	0.426	0.389	0.392	0.384	0.404	0.406	5.68
63)	I	*d5-Chlorobenzene	----------ISTD----------							
64)	TMP	Chlorobenzene	1.201	1.146	1.014	1.044	1.033	0.999	1.073✓	7.59
65)	TM	1,1,1,2-Tetrachloro	0.412	0.404	0.341	0.369	0.377	0.372	0.379	6.75
66)	TMC	Ethylbenzene	1.811	1.639	1.465	1.503	1.504	1.403	1.554	9.51✓
67)	TM	m,p-Xylene	0.739	0.692	0.606	0.599	0.578	0.527	0.623	12.52
68)	TM	o-Xylene	0.673	0.675	0.566	0.573	0.581	0.549	0.603	9.33
69)	TM	Styrene	1.185	1.161	1.012	1.020	1.011	0.980	1.061	8.28
70)	TMP	Bromoform	0.311	0.313	0.301	0.319	0.356	0.352	0.325✓	7.09
71)	TM	Isopropylbenzene	1.853	1.728	1.517	1.539	1.540	1.457	1.606	9.46
72)	TMP	1,1,2,2-Tetrachloro	0.491	0.457	0.414	0.402	0.405	0.386	0.426✓	9.39
73)	S	Bromofluorobenzene	0.651	0.487	0.471	0.461	0.468	0.452	0.498	15.17
74)	TM	1,2,3-Trichloroprop	0.349	0.322	0.315	0.291	0.298	0.282	0.309	7.88
75)	TM	trans-1,4-Dichloro-	0.143	0.135	0.125	0.119	0.121	0.114	0.126	8.59
76)	TM	n-Propylbenzene	2.144	1.962	1.720	1.708	1.660	1.458	1.775	13.63
77)	TM	Bromobenzene	0.497	0.484	0.435	0.431	0.445	0.415	0.451	7.16
78)	TM	1,3,5-Trimethylbenz	1.463	1.381	1.235	1.228	1.244	1.166	1.286	8.69
79)	TM	2-Chlorotoluene	1.405	1.234	1.142	1.145	1.136	1.084	1.191	9.71
80)	TM	4-Chlorotoluene	1.232	1.223	1.026	1.018	1.074	1.009	1.097	9.44
81)	TM	tert-butylbenzene	1.250	1.224	1.074	1.066	1.110	1.020	1.124	8.22
82)	TM	1,2,4-Trimethylbenz	1.490	1.488	1.270	1.201	1.273	1.189	1.319	10.35
83)	T	Diisobutylketone	1.119	1.002	0.874	0.884	0.860	0.775	0.919	13.27
84)	TM	sec-butylbenzene	2.001	1.924	1.692	1.656	1.702	1.645	1.770	8.61
85)	TM	p-isopropyltoluene	1.539	1.540	1.332	1.325	1.349	1.289	1.396	8.10

(#) = Out of Range ### Number of calibration levels exceeded format ###

Response Factor Report Instr_C

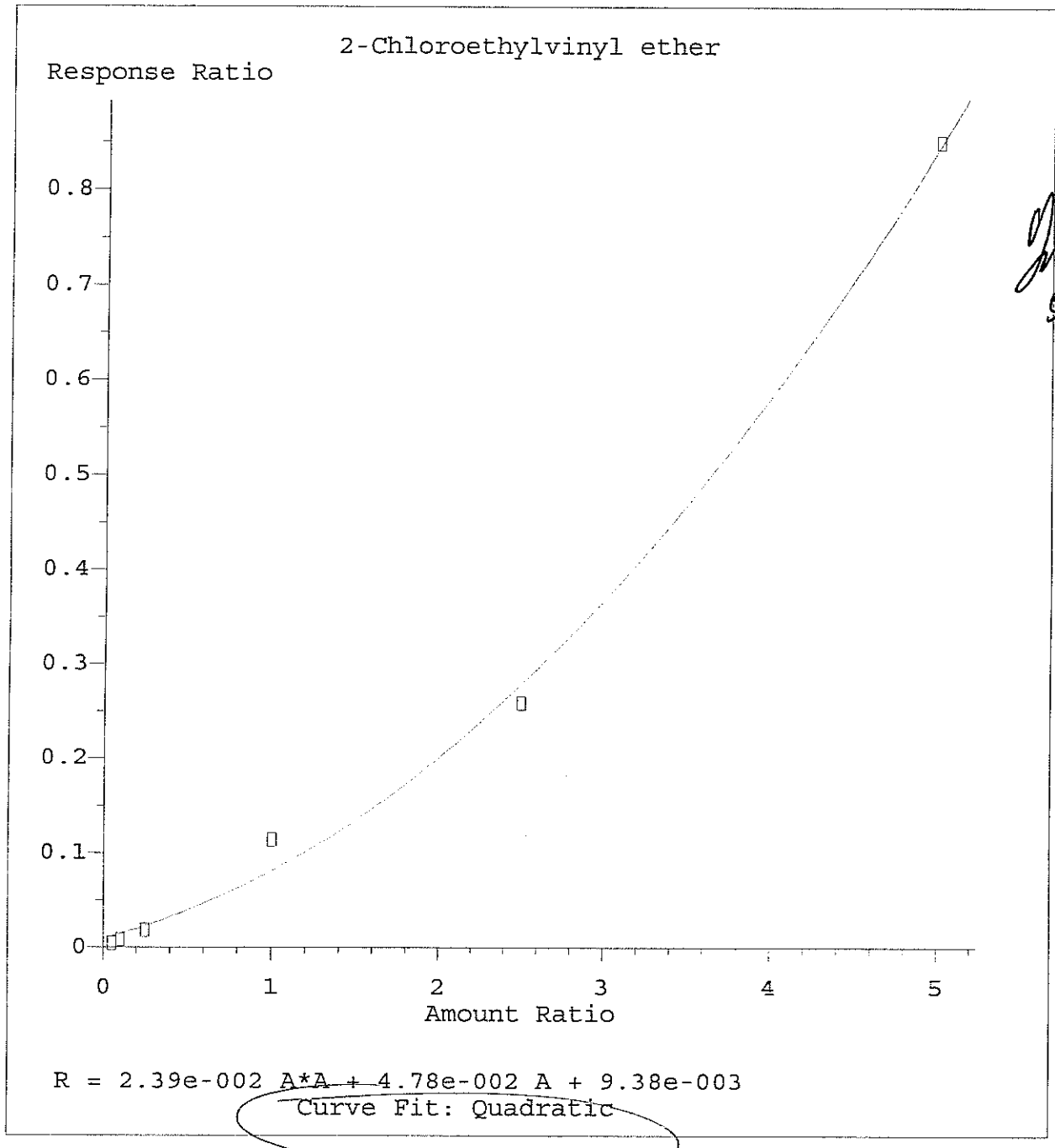
Method : D:\HPCHEM\1\METHODS\V805226C.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Mon May 22 14:18:09 2006
Response via : Initial Calibration

Calibration Files
1 =C59023SI.D 2 =C59024SI.D 5 =C59025SI.D
20 =C59026SI.D 50 =C59027SI.D 100 =C59028SI.D

		Compound	1	2	5	20	50	100	Avg	%RSD
86)	TM	1,3-Dichlorobenzene	0.855	0.804	0.716	0.714	0.730	0.703	0.754	8.16
87)	I	1,4-Dichlorobenzene-d	----------	----------	----------	ISTD	----------	----------	----------	----------
88)	TM	1,4-Dichlorobenzene	1.703	1.663	1.579	1.552	1.556	1.498	1.592	4.83
89)	TM	n-butylbenzene	3.058	2.767	2.528	2.587	2.549	2.514	2.667	7.97
90)	TM	1,2-Dichlorobenzene	1.653	1.474	1.412	1.446	1.403	1.377	1.461	6.86
91)	T	NO COMPOUND							0.000#	-1.00
92)	TM	1,2-Dibromo-3-chlor	0.136	0.168	0.151	0.141	0.150	0.149	0.149	7.26
93)	T	NO COMPOUND							0.000#	-1.00
94)	TM	1,2,4-Trichlorobenz	1.393	1.200	1.145	1.179	1.194	1.216	1.221	7.19
95)	TM	Hexachlorobutadiene	0.679	0.595	0.512	0.581	0.609	0.621	0.599	9.09
96)	TM	Naphthalene	2.681	2.304	2.234	2.153	2.287	2.234	2.315	8.07
97)	TM	1,2,3-Trichlorobenz	1.165	1.067	1.048	1.036	1.101	1.097	1.086	4.29
98)	T	1,3,5-Trichlorobenz	1.493	1.384	1.256	1.305	1.341	1.372	1.359	5.95

(#) = Out of Range ### Number of calibration levels exceeded format ###

MT 5/23/06



5.23.06

Method Name: D:\HPCHEM\1\METHODS\V805226C.M
Calibration Table Last Updated: Mon May 22 14:13:52 2006

Quantitation Report (QT Reviewed)

Data File : D:\HPCHEM\1\DATA\052206C\C59022SI.D Vial: 2
Acq On : 22 May 2006 10:30 am Operator:
Sample : 0.2 PPB Inst : Instr_C
Misc : 5000 Multiplr: 1.00
Integration Params: rteint.p
Quant Time: May 22 14:13 2006 Quant Results File: V805226C.RES

Quant Method : D:\HPCHEM\1\METHODS\V805226C.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Wed Apr 19 15:41:22 2006
Response via : Initial Calibration
DataAcq Meth : V805226C

MT 5/23/06

5-23-06

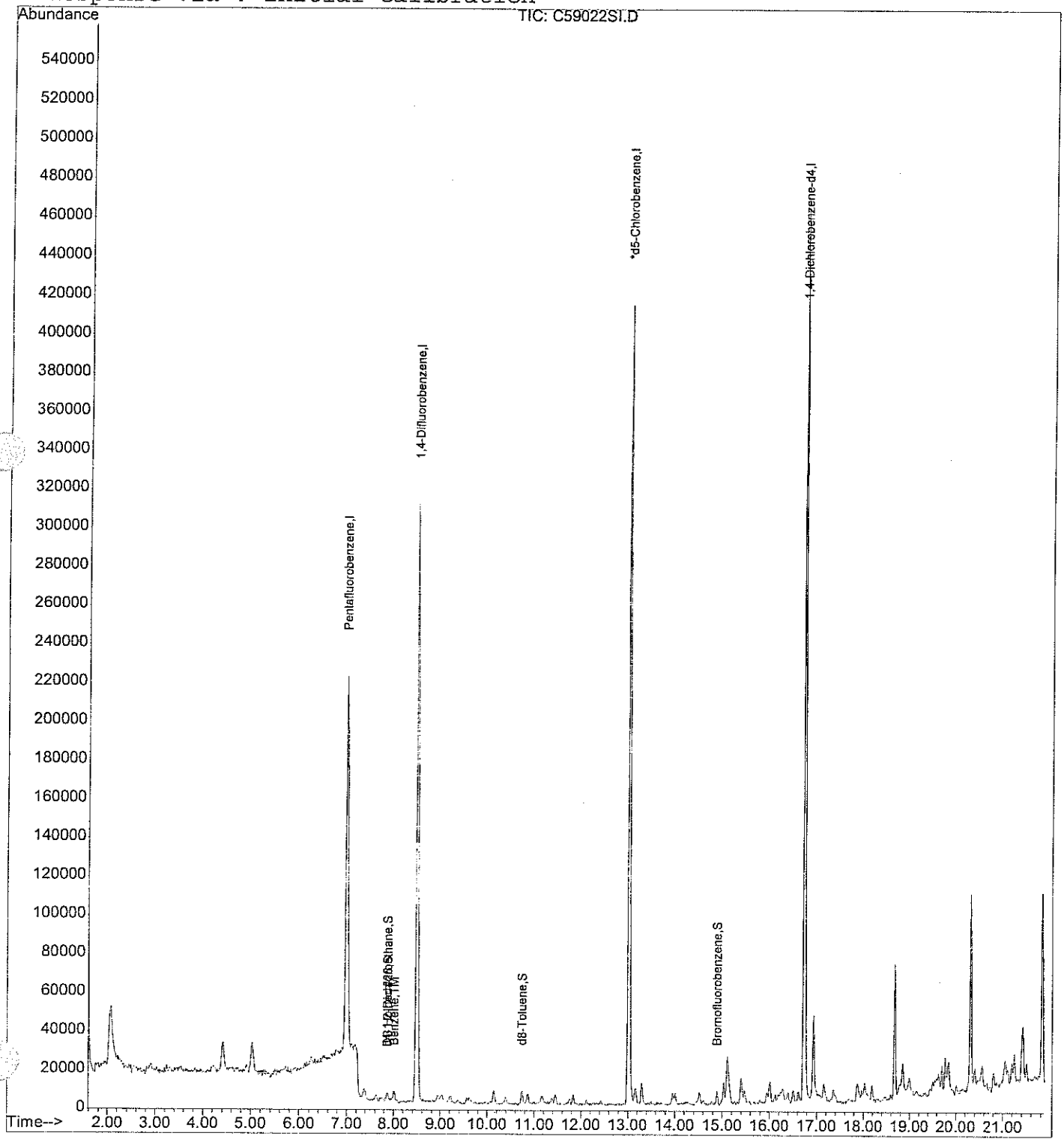
Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	6.99	168	211683	20.00	ug/L	-0.02
39) 1,4-Difluorobenzene	8.48	114	406327 ✓	20.00	ug/L	-0.01
63) *d5-Chlorobenzene	13.02	117	364197 ✓	20.00	ug/L	0.00
87) 1,4-Dichlorobenzene-d4	16.74	152	183801	20.00	ug/L	0.00
System Monitoring Compounds						
28) DB Holder #28	7.87	65	3766	0.45	ug/L	0.00
Spiked Amount 20.000	Range 70 - 130		Recovery = 2.25%#			
36) d4-1,2-Dichloroethane	7.87	65	2645	0.31	ug/L	0.00
Spiked Amount 20.000	Range 70 - 130		Recovery = 1.55%#			
54) d8-Toluene	10.74	98	6185	0.27	ug/L	0.00
Spiked Amount 20.000	Range 70 - 130		Recovery = 1.35%#			
73) Bromofluorobenzene	14.90	95	2966	0.32	ug/L	0.00
Spiked Amount 20.000	Range 70 - 130		Recovery = 1.60%#			
Target Compounds						Qvalue
38) Benzene	8.00	78	5476m	0.32	ug/L	

(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : D:\HPCHEM\1\DATA\052206C\C59022SI.D Vial: 2
Acq On : 22 May 2006 10:30 am Operator:
Sample : 0.2 PPB Inst : Instr_C
Misc : 5000 Multiplr: 1.00
MS Integration Params: rteint.p
Quant Time: May 22 14:13 2006 Quant Results File: V805226C.RES

Method : D:\HPCHEM\1\METHODS\V805226C.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Mon May 22 14:18:09 2006
Response via : Initial Calibration

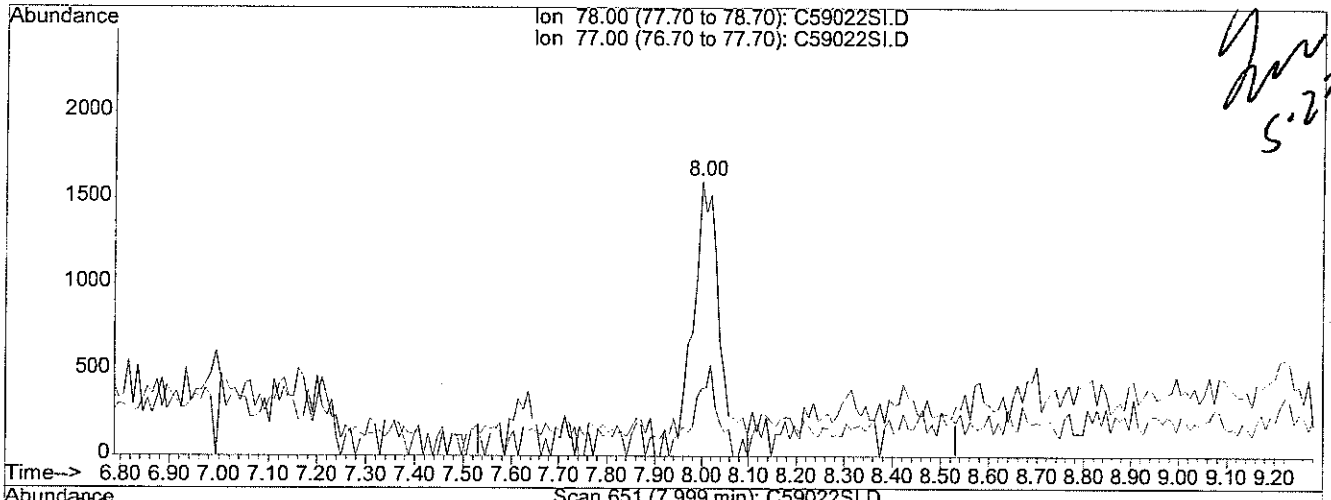


Quantitation Report (Qedit)

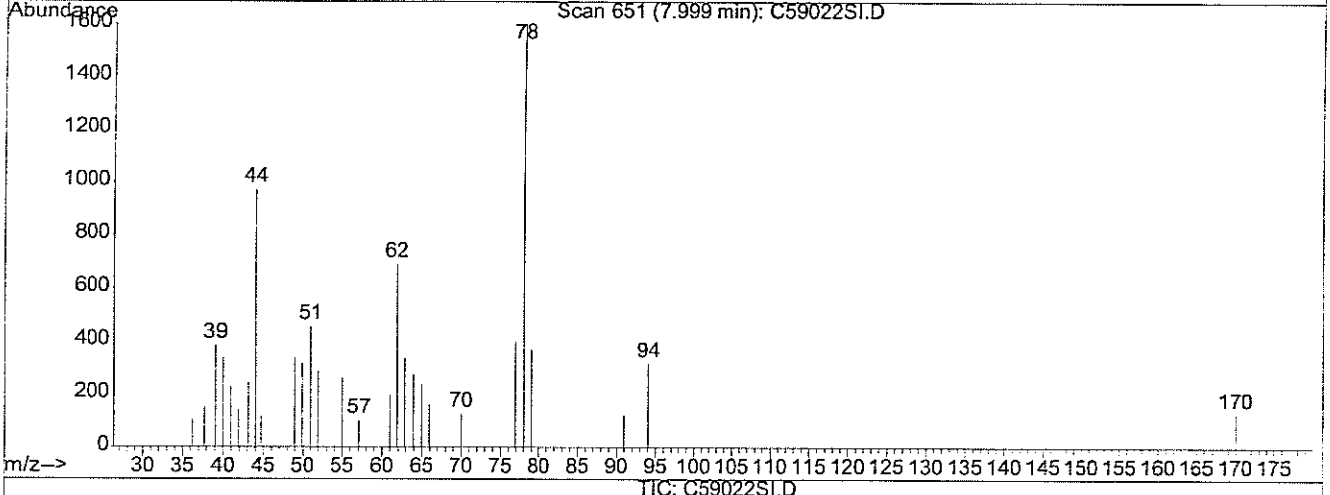
Data File : D:\HPCHEM\1\DATA\052206C\C59022SI.D Vial: 2
Acq On : 22 May 2006 10:30 am Operator:
Sample : 0.2 PPB Inst : Instr_C
Misc : 5000 Multiplr: 1.00
MS Integration Params: rteint.p
Quant Time: May 22 14:04 2006 Quant Results File: temp.res

Method : D:\HPCHEM\1\METHODS\V805226C.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Mon May 22 14:09:16 2006
Response via : Multiple Level Calibration

MT 5/23/06
extra area



5.25.06



TIC: C59022SI.D

(38) Benzene (TM)

8.00min 0.40ug/L

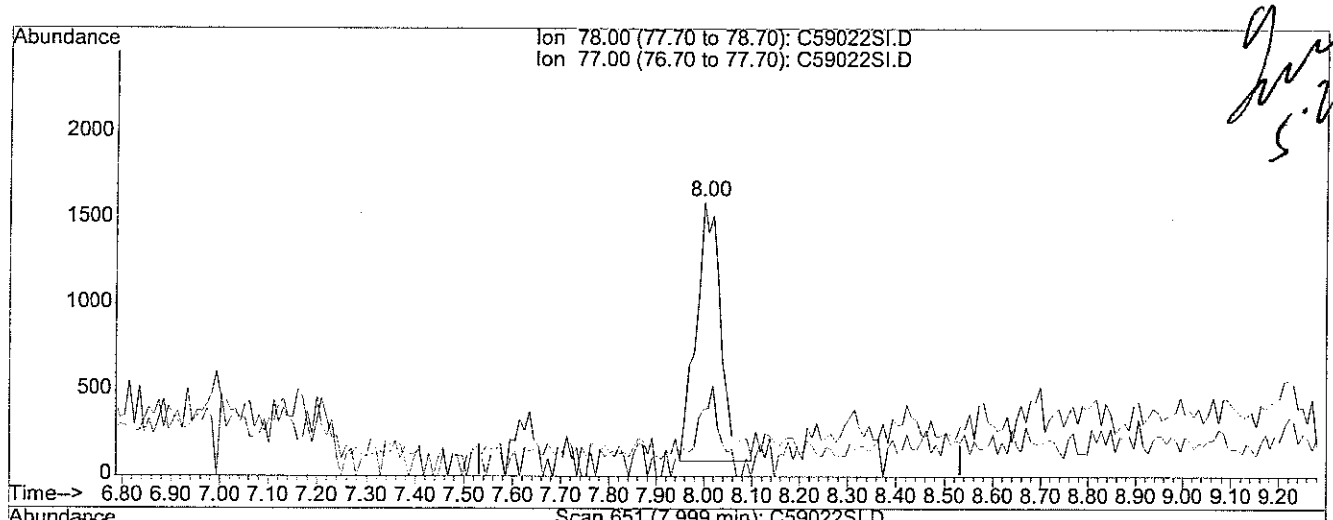
response 6862

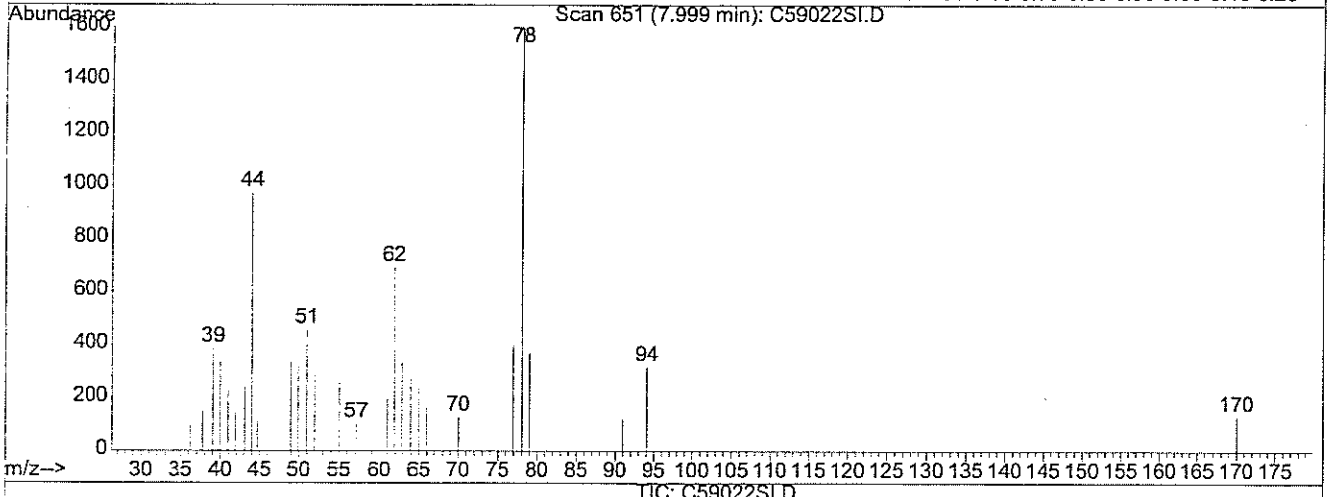
Ion	Exp%	Act%
78.00	100	100
77.00	24.00	24.80
0.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (Qedit)

Data File : D:\HPCHEM\1\DATA\052206C\C59022SI.D Vial: 2
Acq On : 22 May 2006 10:30 am Operator:
Sample : 0.2 PPB Inst : Instr_C
Misc : 5000 Multiplr: 1.00
MS Integration Params: rteint.p
Quant Time: May 22 14:13 2006 Quant Results File: temp.res

Method : D:\HPCHEM\1\METHODS\V805226C.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Mon May 22 14:09:16 2006 MT 5/23/06
Response via : Multiple Level Calibration





TIC: C59022SI.D

(38) Benzene (TM)

8.00min 0.32ug/L m

response 5476

Ion	Exp%	Act%
78.00	100	100
77.00	24.00	24.80
0.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (QT Reviewed)

Data File : D:\HPCHEM\1\DATA\052206C\C59023SI.D Vial: 3
Acq On : 22 May 2006 11:24 am Operator:
Sample : 1 PPB Inst : Instr_C
Misc : 5000 Multiplr: 1.00
Integration Params: rteint.p
Quant Time: May 22 14:11 2006 Quant Results File: V805226C.RES

Quant Method : D:\HPCHEM\1\METHODS\V805226C.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Wed Apr 19 15:41:22 2006
Response via : Initial Calibration
DataAcq Meth : V805226C

MT 5/23/06

5-23-06

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.00	168	250423	20.00	ug/L	-0.02
39) 1,4-Difluorobenzene	8.49	114	437021 ✓	20.00	ug/L	0.00
63) *d5-Chlorobenzene	13.02	117	379373 ✓	20.00	ug/L	0.00
87) 1,4-Dichlorobenzene-d4	16.74	152	185550	20.00	ug/L	0.00
System Monitoring Compounds						
28) DB Holder #28	7.88	65	12403	1.25	ug/L	0.00
Spiked Amount 20.000	Range 70 - 130		Recovery = 6.25%#			
36) d4-1,2-Dichloroethane	7.88	65	12403	1.25	ug/L	0.00
Spiked Amount 20.000	Range 70 - 130		Recovery = 6.25%#			
54) d8-Toluene	10.75	98	27872	1.14	ug/L	0.00
Spiked Amount 20.000	Range 70 - 130		Recovery = 5.70%#			
73) Bromofluorobenzene	14.89	95	12342	1.28	ug/L	0.00
Spiked Amount 20.000	Range 70 - 130		Recovery = 6.40%#			
Target Compounds						Qvalue
2) Dichlorodifluoromethane	1.70	85	8847	1.05	ug/L	95
3) Chloromethane	1.95	50	8156m	0.96	ug/L	
4) Vinyl Chloride	2.03	62	9749m	1.12	ug/L	
5) Bromomethane	2.49	94	4042	0.90	ug/L #	75
6) Chloroethane	2.56	64	5742m	1.35	ug/L	
7) t-Butyl alcohol (TBA)	3.86	59	5503	8.46	ug/L #	89
8) Trichlorofluoromethane	2.88	101	18649	1.19	ug/L	99
9) Diethyl ether	3.26	74	2656m	1.08	ug/L	
10) 1,1,2-Trichlorotrifluoroet	3.46	101	7665m	1.25	ug/L	
11) Acrolein	3.42	56	6901	12.76	ug/L	93
12) Acetone	3.52	43	21552	9.92	ug/L	100
13) 1,1-Dichloroethene	3.66	96	6627	1.49	ug/L #	57
14) Methyl iodide	4.10	142	8768	1.27	ug/L	94
15) Di-isopropyl ether (DIPE)	5.50	45	25837	1.18	ug/L	97
16) Methylene Chloride	4.43	84	9441	1.94	ug/L #	79
17) Carbon Disulfide	4.39	76	16598	1.16	ug/L	100
18) Acrylonitrile	4.63	53	2956	1.44	ug/L #	84
19) Methyl-tert-butyl ether (M	4.73	73	21636	1.29	ug/L #	93
20) trans-1,2-Dichloroethene	4.90	61	13356	1.22	ug/L	96
21) 1,1-Dichloroethane	5.59	63	13991	1.12	ug/L	89
22) Vinyl acetate	5.65	43	17572	1.69	ug/L	98
23) Methyl ethyl ketone	6.31	72	4174	7.86	ug/L	66
24) Ethyl t-butyl ether (ETBE)	6.14	59	26713	1.30	ug/L	94
25) 2,2-Dichloropropane	6.46	77	14594m	1.52	ug/L	
26) cis-1,2-Dichloroethene	6.54	96	7779	1.10	ug/L #	74
27) t-Amyl methyl ether (TAME)	7.88	73	19141	1.33	ug/L #	92

(#) = qualifier out of range (m) = manual integration

Quantitation Report (QT Reviewed)

Data File : D:\HPCHEM\1\DATA\052206C\C59023SI.D Vial: 3
Acq On : 22 May 2006 11:24 am Operator:
Sample : 1 PPB Inst : Instr_C
Misc : 5000 Multiplr: 1.00
Integration Params: rteint.p
Quant Time: May 22 14:11 2006 Quant Results File: V805226C.RES

Quant Method : D:\HPCHEM\1\METHODS\V805226C.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Wed Apr 19 15:41:22 2006
Response via : Initial Calibration
DataAcq Meth : V805226C

	Compound	R.T.	QIon	Response	Conc Unit	Qvalue
29)	Chloroform	6.78	83	14469	1.23 ug/L	94
30)	Bromochloromethane	7.02	128	4612	1.23 ug/L	90
31)	Tetrahydrofuran	7.12	71	1012	1.76 ug/L #	64
32)	1,1,1-Trichloroethane	7.38	97	12840	1.20 ug/L	89
34)	1,1-Dichloropropene	7.63	75	10933	1.28 ug/L	91
35)	Carbon Tetrachloride	7.76	119	9649	1.09 ug/L	92
37)	1,2-Dichloroethane	8.02	62	13754	1.18 ug/L	90
38)	Benzene	8.02	78	26348	1.30 ug/L	92
40)	Trichloroethene	8.95	95	9306	1.21 ug/L	92
41)	1,2-Dichloropropane	9.21	63	8203	1.13 ug/L	98
42)	Methylmethacrylate	9.37	69	5136	1.15 ug/L	93
43)	Methyl acetate	4.23	43	6961	1.05 ug/L #	72
44)	Cyclohexane	7.39	56	14691	1.26 ug/L	95
45)	Bromodichloromethane	9.56	83	10030	1.06 ug/L	98
46)	Dibromomethane	9.62	93	6356	1.15 ug/L	95
47)	1,4-Dioxane	9.63	88	2281	33.37 ug/L #	59
48)	Methylcyclohexane	9.03	83	12147	1.30 ug/L	96
50)	2-Chloroethylvinyl ether	10.08	63	2025	0.95 ug/L #	84
51)	2-Hexanone	11.44	58	15950	6.62 ug/L	99
52)	Methyl isobutyl ketone	10.14	58	16054	6.46 ug/L	100
53)	cis-1,3-Dichloropropene	10.39	75	11312	1.05 ug/L	93
55)	Toluene	10.87	92	17640	1.22 ug/L	100
56)	trans-1,3-Dichloropropene	11.15	75	11409	1.11 ug/L	98
58)	1,1,2-Trichloroethane	11.38	83	5613	1.18 ug/L	95
59)	1,3-Dichloropropane	11.76	76	11268	1.21 ug/L	96
60)	Tetrachloroethene	11.84	166	10390	1.11 ug/L	95
61)	Dibromochloromethane	12.13	129	8735	1.09 ug/L	96
62)	1,2-Dibromoethane	12.42	107	9656	1.20 ug/L	97
64)	Chlorobenzene	13.07	112	22782	1.19 ug/L	94
65)	1,1,1,2-Tetrachloroethane	13.15	131	7815	1.15 ug/L	94
66)	Ethylbenzene	13.17	91	34345	1.24 ug/L	94
67)	m,p-Xylene	13.29	106	28043	2.53 ug/L	94
68)	o-Xylene	13.96	106	12762	1.19 ug/L	93
69)	Styrene	14.01	104	22476	1.19 ug/L	91
70)	Bromoform	14.49	173	5892	1.00 ug/L	99
71)	Isopropylbenzene	14.52	105	35158	1.24 ug/L	98
72)	1,1,2,2-Tetrachloroethane	14.78	83	9322	1.25 ug/L	94
74)	1,2,3-Trichloropropane	15.00	75	6625	1.20 ug/L #	89
75)	trans-1,4-Dichloro-2-buten	15.14	89	2718	1.18 ug/L	93
76)	n-Propylbenzene	15.15	91	40676	1.28 ug/L	98
77)	Bromobenzene	15.17	156	9423	1.15 ug/L	85

(#) = qualifier out of range (m) = manual integration

Quantitation Report (QT Reviewed)

Data File : D:\HPCHEM\1\DATA\052206C\C59023SI.D Vial: 3
Acq On : 22 May 2006 11:24 am Operator:
Sample : 1 PPB Inst : Instr_C
Misc : 5000 Multiplr: 1.00
Integration Params: rteint.p
Quant Time: May 22 14:11 2006 Quant Results File: V805226C.RES

Quant Method : D:\HPCHEM\1\METHODS\V805226C.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Wed Apr 19 15:41:22 2006
Response via : Initial Calibration
DataAcq Meth : V805226C

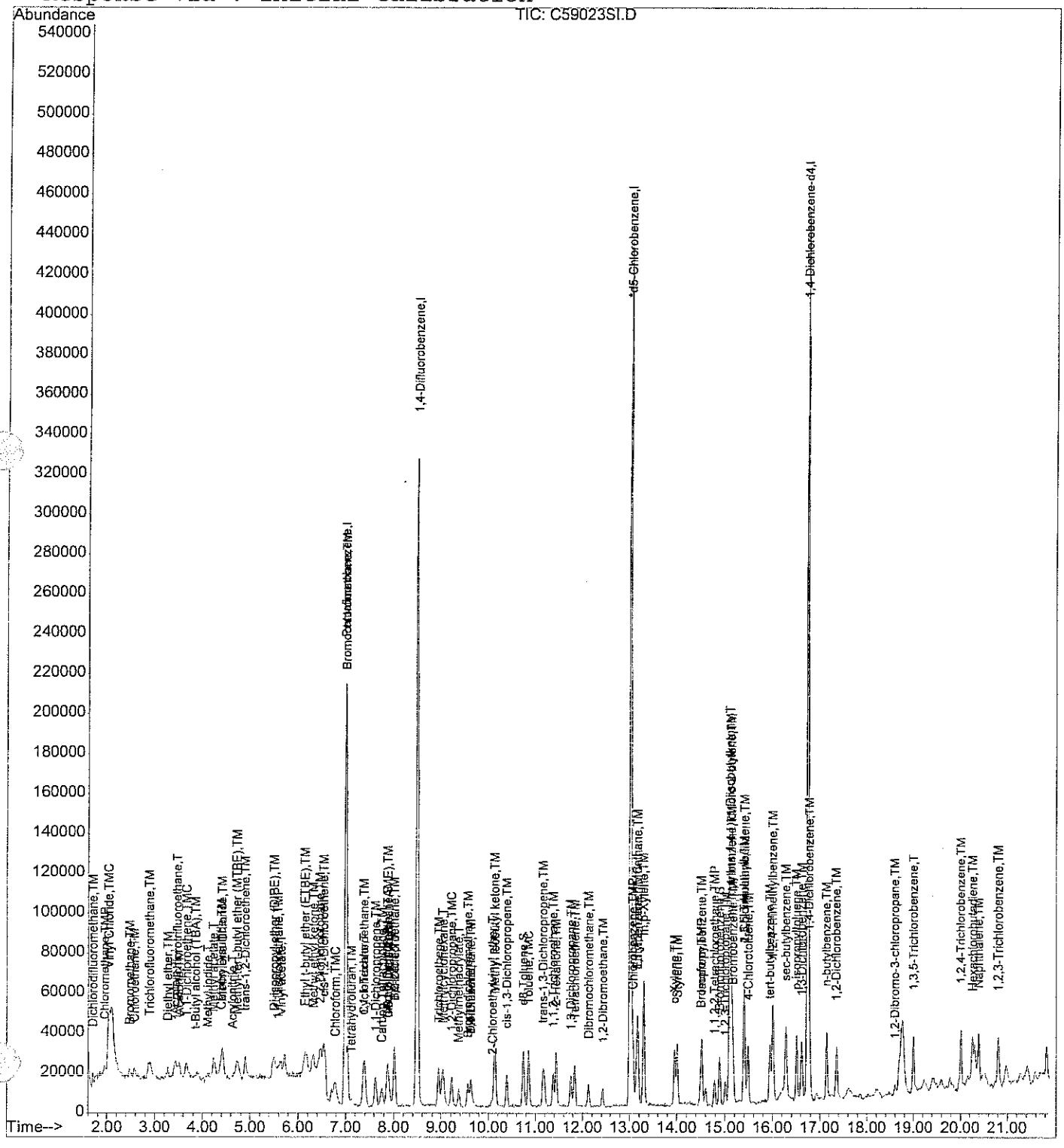
	Compound	R.T.	QIon	Response	Conc Unit		Qvalue
78)	1,3,5-Trimethylbenzene	15.40	105	27757	1.18 ug/L		97
79)	2-Chlorotoluene	15.40	91	26657	1.21 ug/L		97
80)	4-Chlorotoluene	15.48	91	23364	1.16 ug/L		100
81)	tert-butylbenzene	15.94	119	23709	1.14 ug/L		99
82)	1,2,4-Trimethylbenzene	16.00	105	28271	1.18 ug/L		97
83)	Diisobutylketone	15.11	57	106163	6.35 ug/L		92
84)	sec-butylbenzene	16.28	105	37952	1.19 ug/L		98
85)	p-isopropyltoluene	16.51	119	29192	1.14 ug/L		98
86)	1,3-Dichlorobenzene	16.62	146	16211	1.17 ug/L		97
88)	1,4-Dichlorobenzene	16.78	146	15804	1.14 ug/L		92
89)	n-butylbenzene	17.15	91	28371	1.22 ug/L		97
90)	1,2-Dichlorobenzene	17.35	146	15337	1.20 ug/L		99
92)	1,2-Dibromo-3-chloropropan	18.60	75	1266	0.93 ug/L	#	82
94)	1,2,4-Trichlorobenzene	20.01	180	12927	1.18 ug/L		100
95)	Hexachlorobutadiene	20.25	225	6296	1.12 ug/L		98
96)	Naphthalene	20.39	128	24876	1.24 ug/L		100
97)	1,2,3-Trichlorobenzene	20.81	180	10809	1.08 ug/L		94
98)	1,3,5-Trichlorobenzene	19.00	180	13854	1.12 ug/L		98

(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : D:\HPCHEM\1\DATA\052206C\C59023SI.D Vial: 3
Acq On : 22 May 2006 11:24 am Operator:
Sample : 1 PPB Inst : Instr_C
Misc : 5000 Multiplr: 1.00
MS Integration Params: rteint.p
Quant Time: May 22 14:11 2006 Quant Results File: V805226C.RES

Method : D:\HPCHEM\1\METHODS\V805226C.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Mon May 22 14:18:09 2006
Response via : Initial Calibration

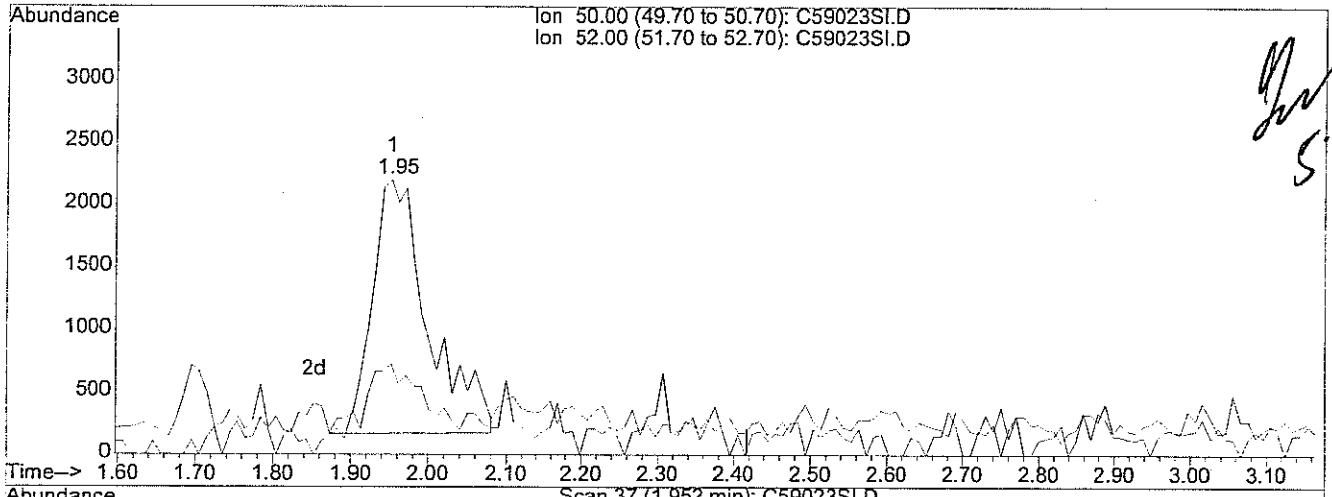


Quantitation Report (Qedit)

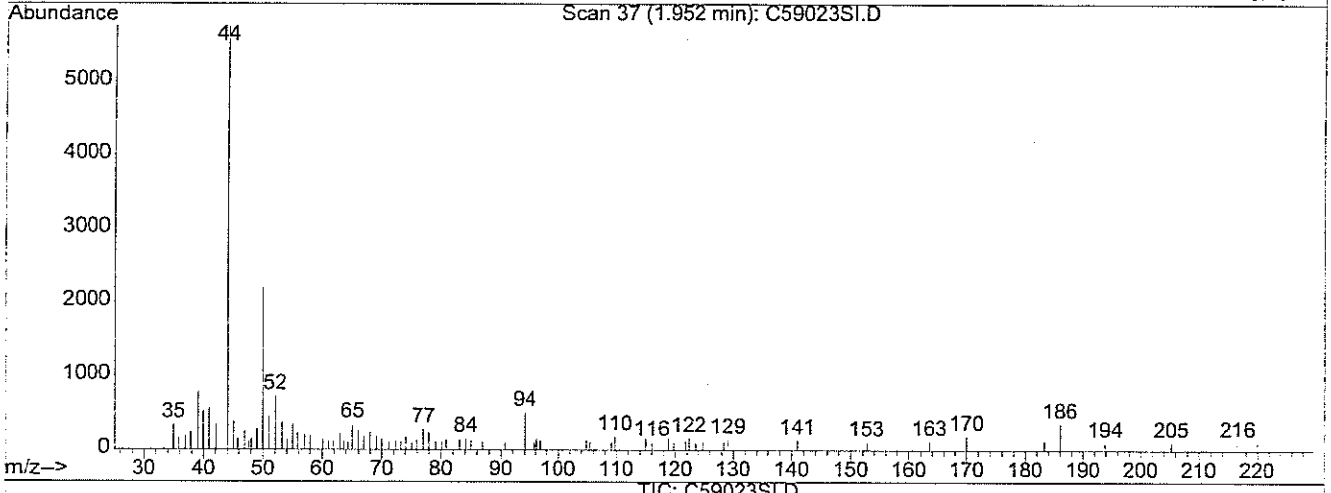
Data File : D:\HPCHEM\1\DATA\052206C\C59023SI.D Vial: 3
Acq On : 22 May 2006 11:24 am Operator:
Sample : 1 PPB Inst : Instr_C
Misc : 5000 Multiplr: 1.00
MS Integration Params: rteint.p
Quant Time: May 22 13:19 2006 Quant Results File: temp.res

Method : D:\HPCHEM\1\METHODS\V805226C.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Mon May 22 14:05:38 2006
Response via : Multiple Level Calibration

MT 5/23/06
extra area



5.23.06



TIC: C59023SI.D

(3) Chloromethane (TMP)

1.95min 1.21ug/L

response 10256

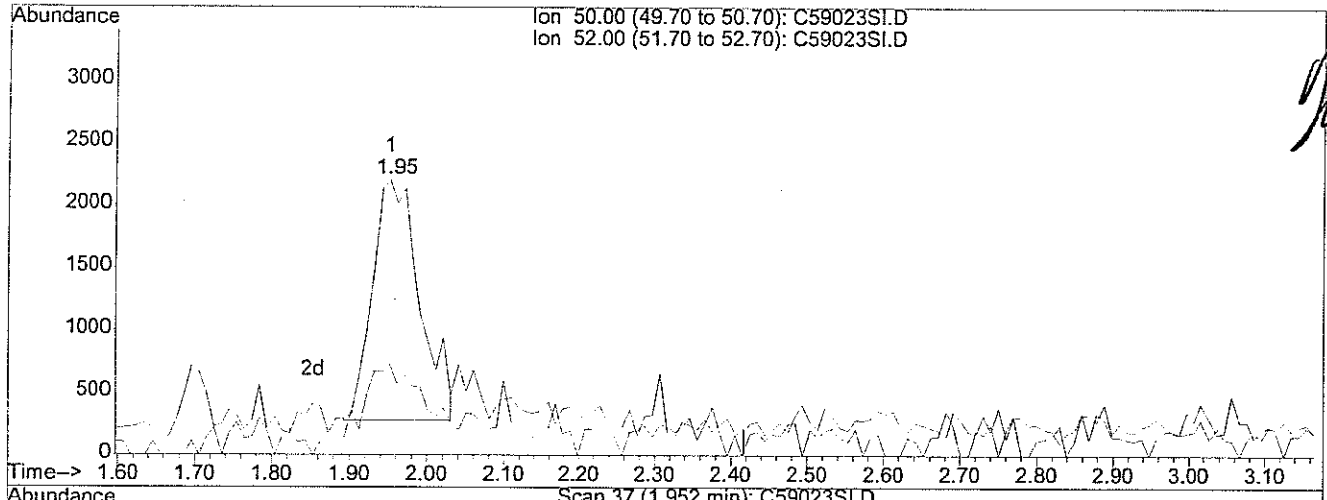
Ion	Exp%	Act%
50.00	100	100
52.00	31.40	28.23
0.00	0.00	0.00
0.00	0.00	0.00

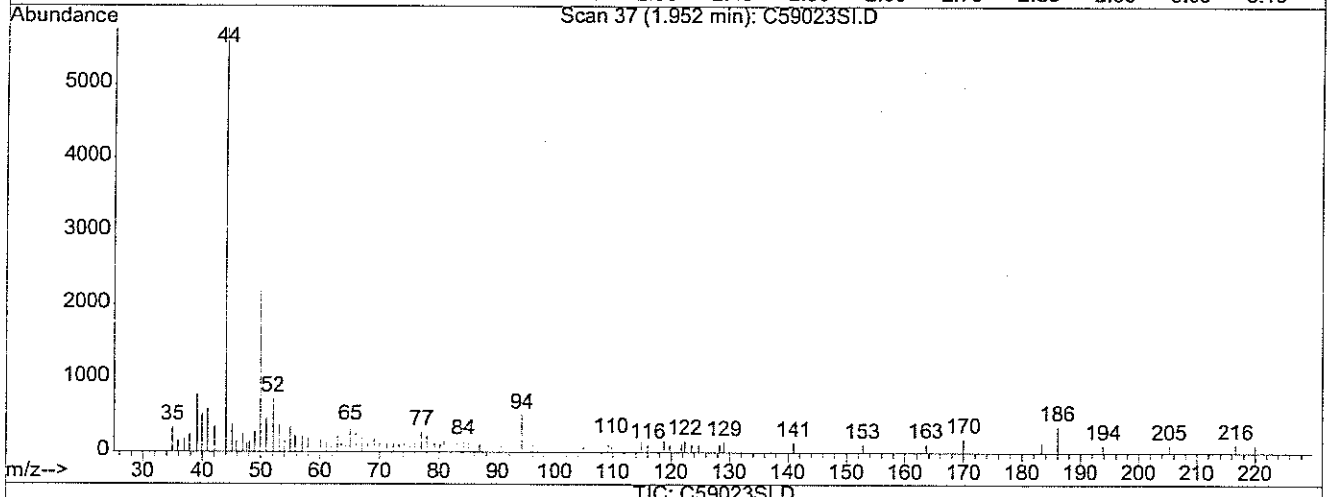
Quantitation Report (Qedit)

Data File : D:\HPCHEM\1\DATA\052206C\C59023SI.D Vial: 3
Acq On : 22 May 2006 11:24 am Operator:
Sample : 1 PPB Inst : Instr_C
Misc : 5000 Multiplr: 1.00
MS Integration Params: rteint.p
Quant Time: May 22 14:07 2006 Quant Results File: temp.res

Method : D:\HPCHEM\1\METHODS\V805226C.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Mon May 22 14:05:38 2006
Response via : Multiple Level Calibration

MT 5/23/06





5-23-06

(3) Chloromethane (TMP)

1.95min 0.96ug/L m

response 8156

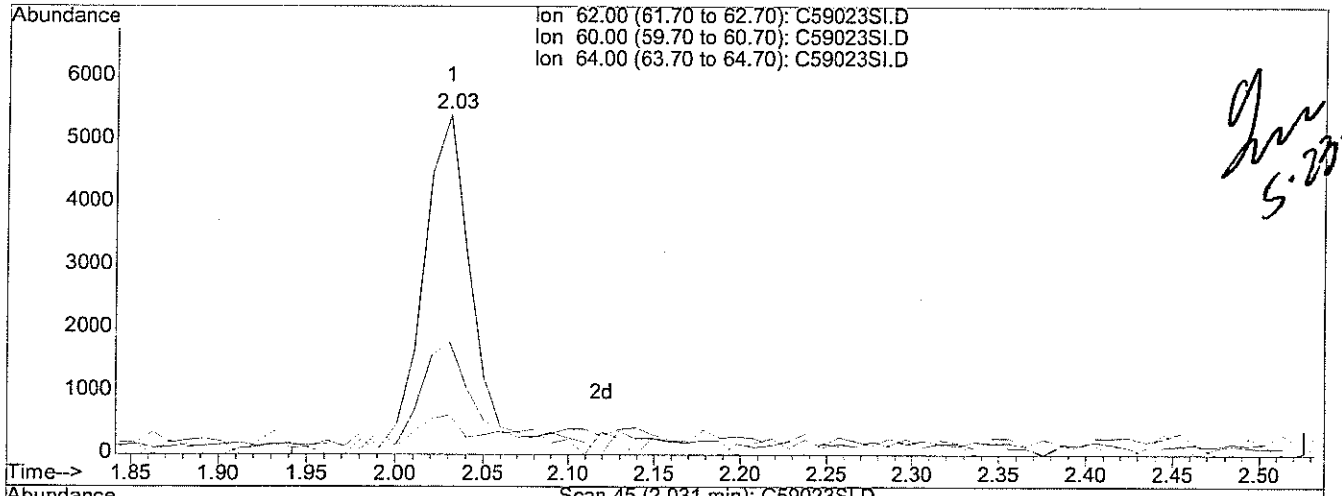
Ion	Exp%	Act%
50.00	100	100
52.00	31.40	33.03
0.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (Qedit)

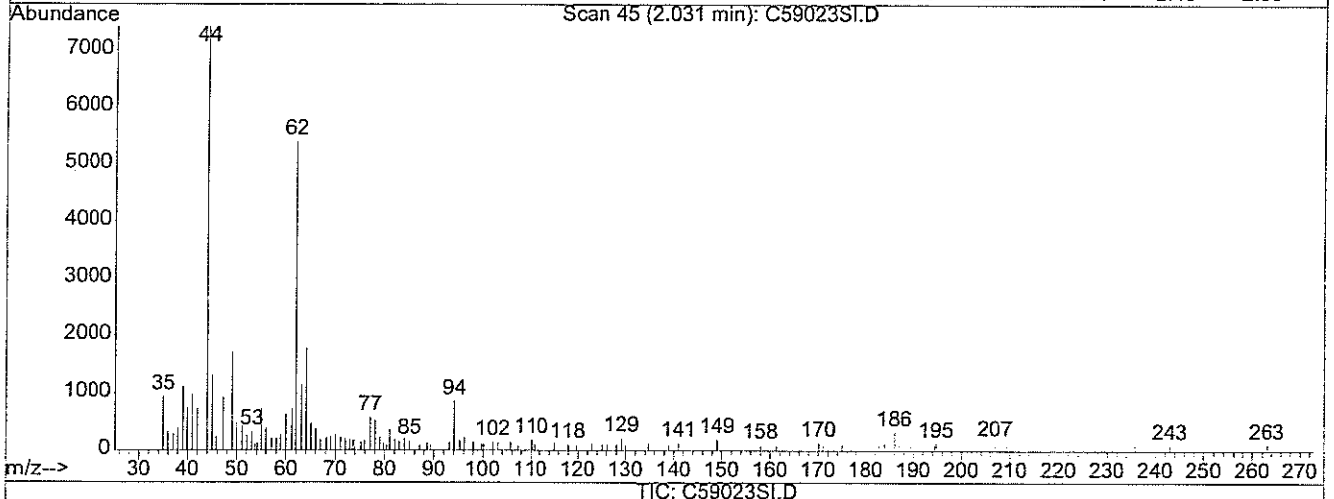
Data File : D:\HPCHEM\1\DATA\052206C\C59023SI.D Vial: 3
Acq On : 22 May 2006 11:24 am Operator:
Sample : 1 PPB Inst : Instr_C
Misc : 5000 Multiplr: 1.00
MS Integration Params: rteint.p
Quant Time: May 22 14:07 2006 Quant Results File: temp.res

Method : D:\HPCHEM\1\METHODS\V805226C.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Mon May 22 14:05:38 2006
Response via : Multiple Level Calibration

MT 5/23/06
extra area



5.23.06



TIC: C59023SI.D

(4) Vinyl Chloride (TMC)

2.03min 1.30ug/L

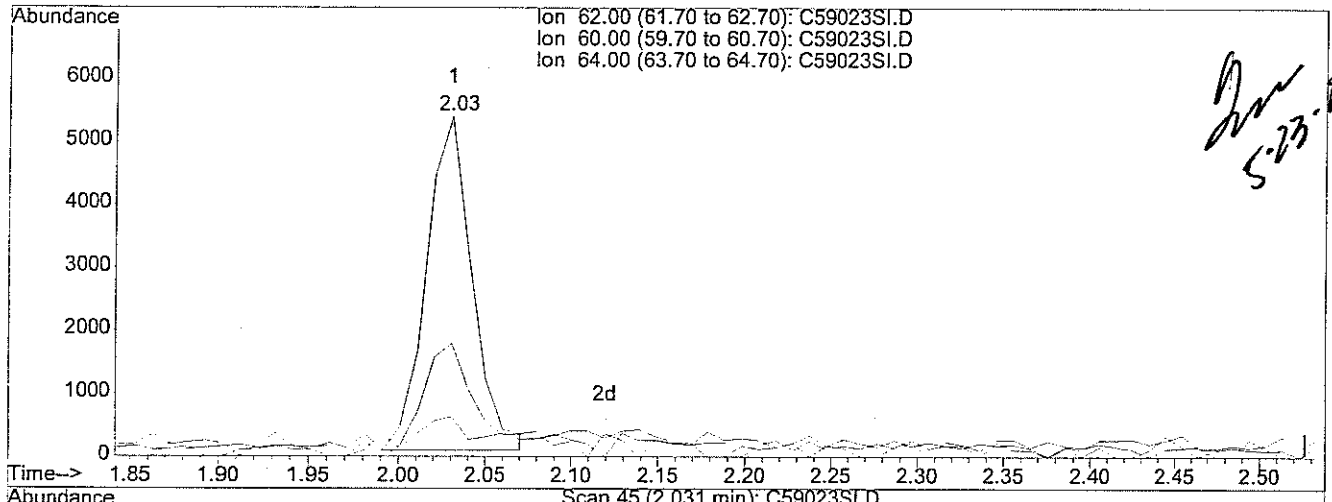
response 11244

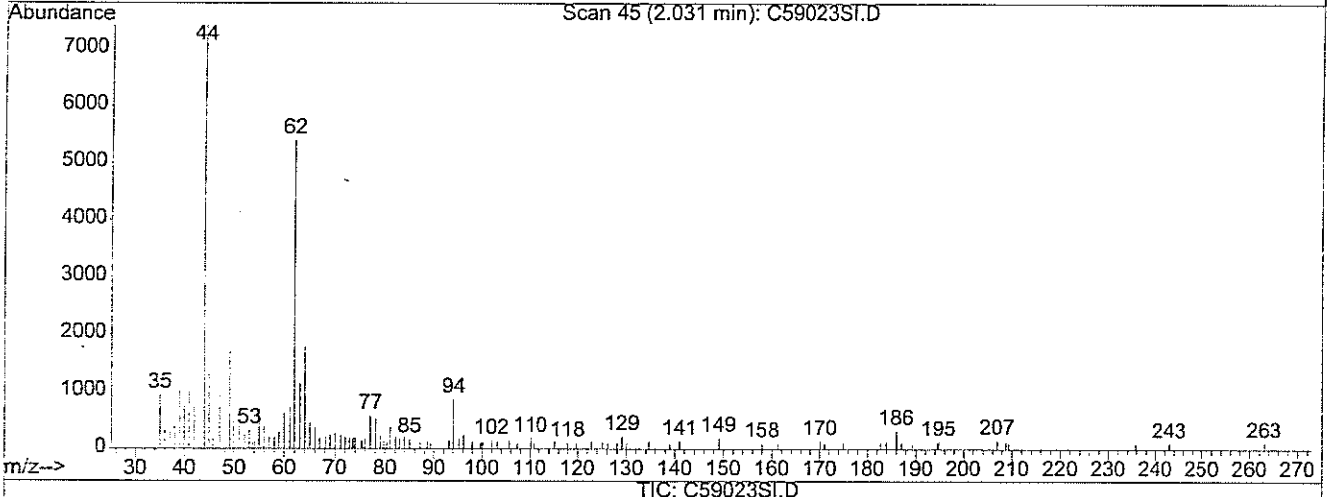
Ion	Exp%	Act%
62.00	100	100
60.00	6.30	8.90
64.00	31.70	29.95
0.00	0.00	0.00

Quantitation Report (Qedit)

Data File : D:\HPCHEM\1\DATA\052206C\C59023SI.D Vial: 3
Acq On : 22 May 2006 11:24 am Operator:
Sample : 1 PPB Inst : Instr_C
Misc : 5000 Multiplr: 1.00
MS Integration Params: rteint.p
Quant Time: May 22 14:07 2006 Quant Results File: temp.res

Method : D:\HPCHEM\1\METHODS\V805226C.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Mon May 22 14:05:38 2006 MT 5/23/06
Response via : Multiple Level Calibration





TIC: C59023SI.D

(4) Vinyl Chloride (TMC)

2.03min 1.12ug/L m

response 9749

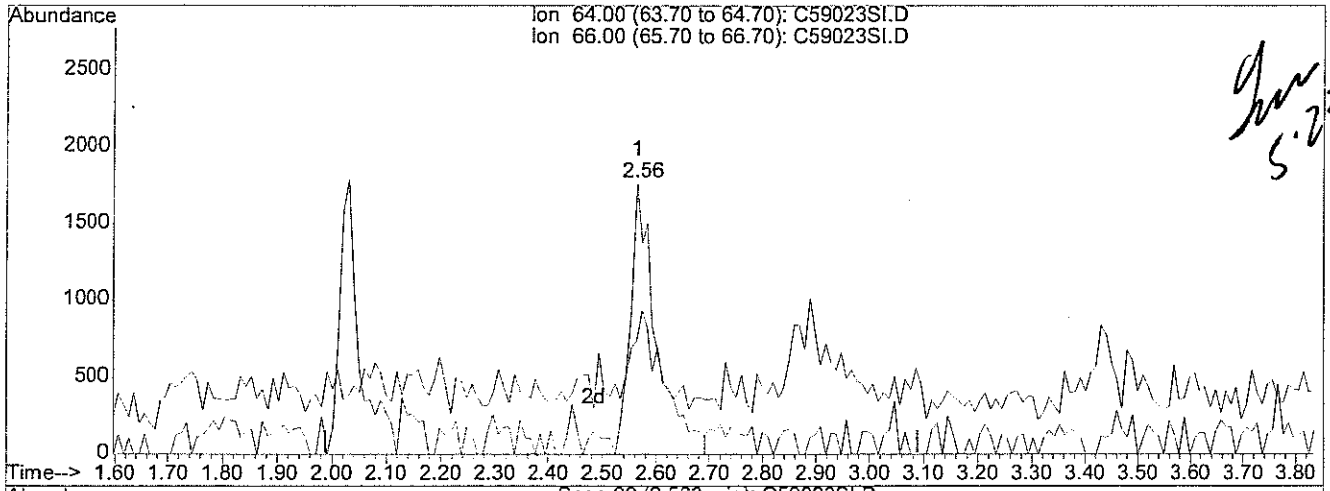
Ion	Exp%	Act%
62.00	100	100
60.00	6.30	11.66
64.00	31.70	33.17
0.00	0.00	0.00

Quantitation Report (Qedit)

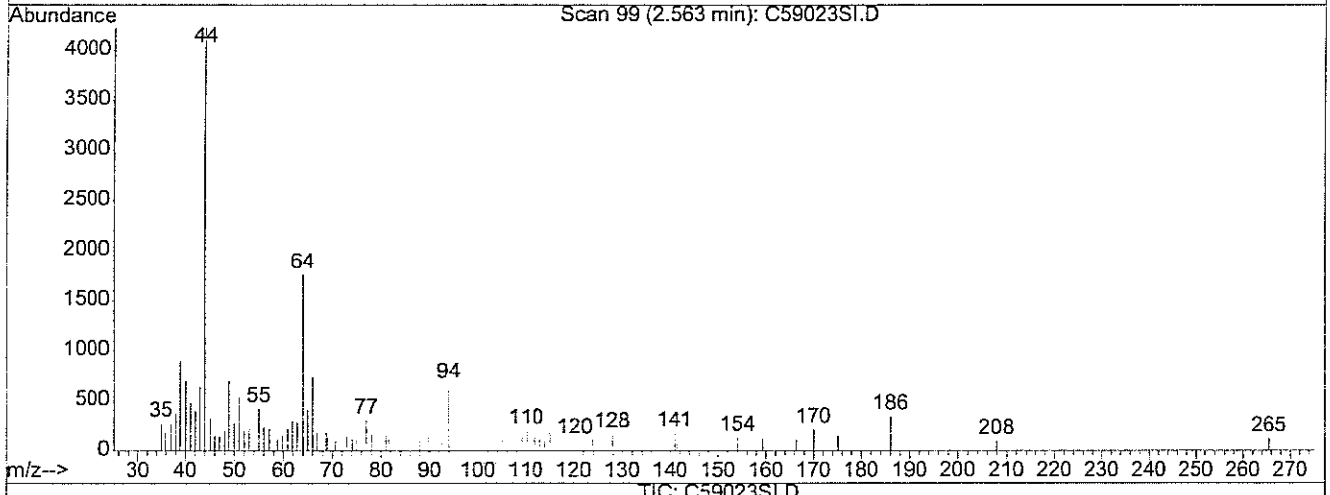
Data File : D:\HPCHEM\1\DATA\052206C\C59023SI.D Vial: 3
Acq On : 22 May 2006 11:24 am Operator:
Sample : 1 PPB Inst : Instr_C
Misc : 5000 Multiplr: 1.00
MS Integration Params: rteint.p
Quant Time: May 22 14:07 2006 Quant Results File: temp.res

Method : D:\HPCHEM\1\METHODS\V805226C.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Mon May 22 14:05:38 2006
Response via : Multiple Level Calibration

MT 5/23/06
extra area



5.23.06



TIC: C59023SI.D

(6) Chloroethane (TM)

2.56min 1.41ug/L

response 5994

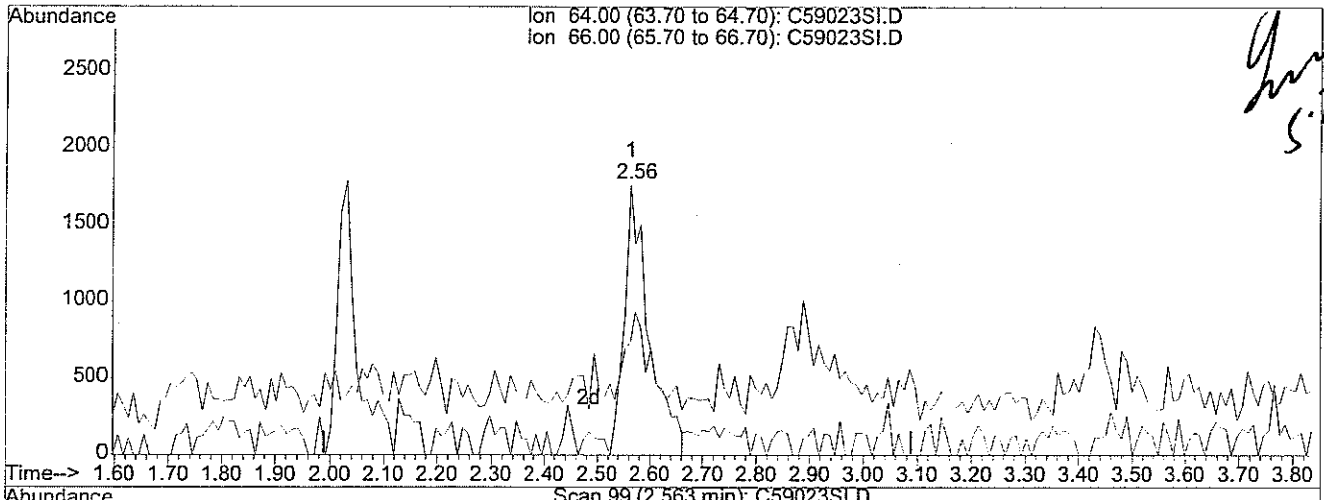
Ion	Exp%	Act%
64.00	100	100
66.00	31.70	21.77
0.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (Qedit)

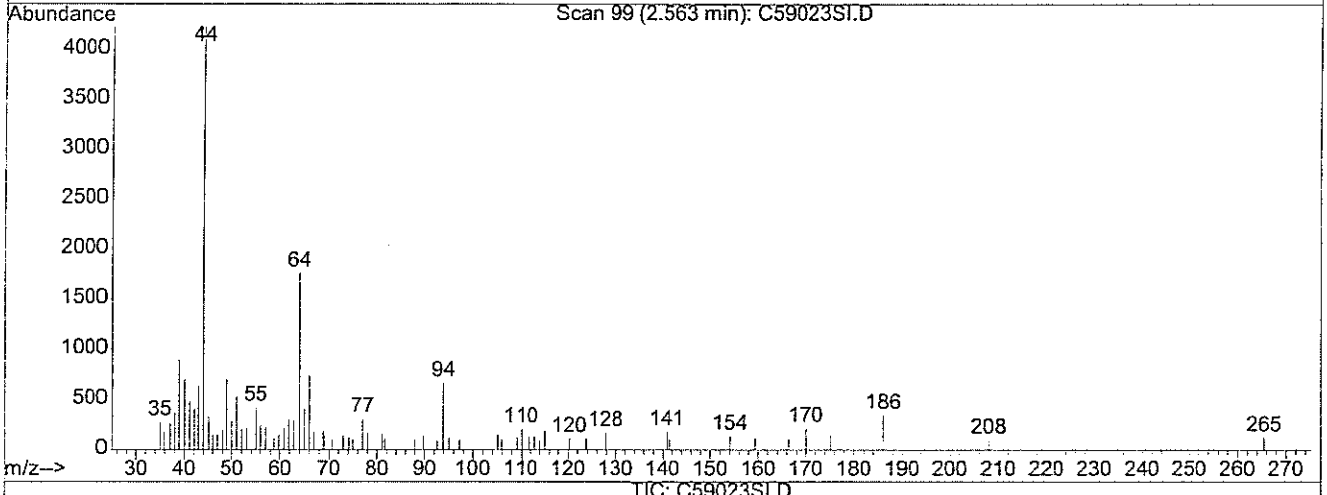
Data File : D:\HPCHEM\1\DATA\052206C\C59023SI.D Vial: 3
Acq On : 22 May 2006 11:24 am Operator:
Sample : 1 PPB Inst : Instr_C
Misc : 5000 Multiplr: 1.00
MS Integration Params: rteint.p
Quant Time: May 22 14:08 2006 Quant Results File: temp.res

Method : D:\HPCHEM\1\METHODS\V805226C.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Mon May 22 14:05:38 2006
Response via : Multiple Level Calibration

MT 5/23/06



5.23.06



TIC: C59023SI.D

(6) Chloroethane (TM)

2.56min 1.35ug/L m

response 5742

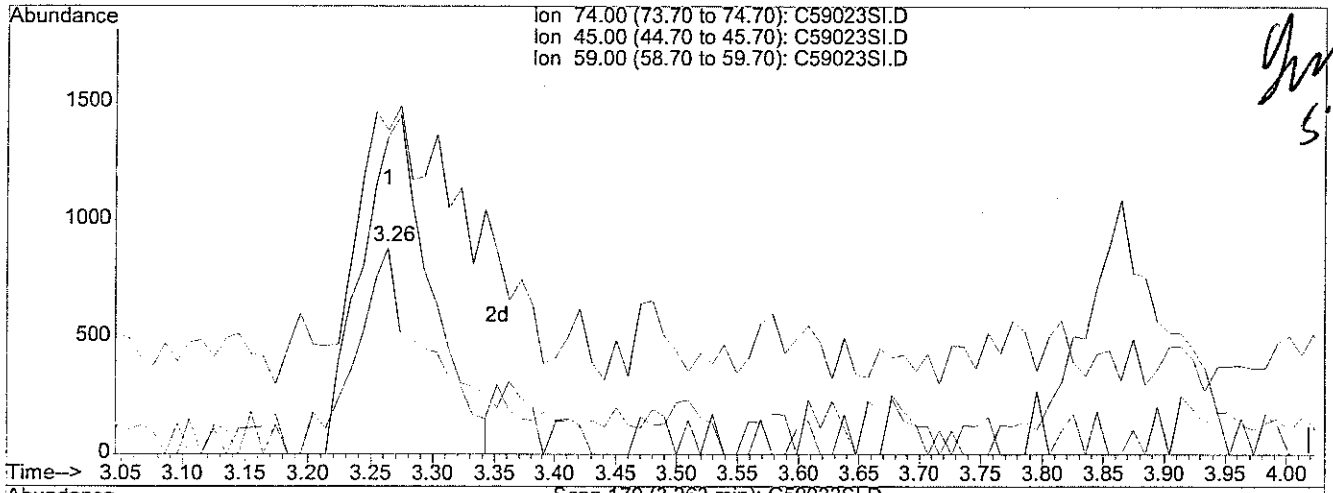
Ion	Exp%	Act%
64.00	100	100
66.00	31.70	42.12
0.00	0.00	0.00
0.00	0.00	0.00

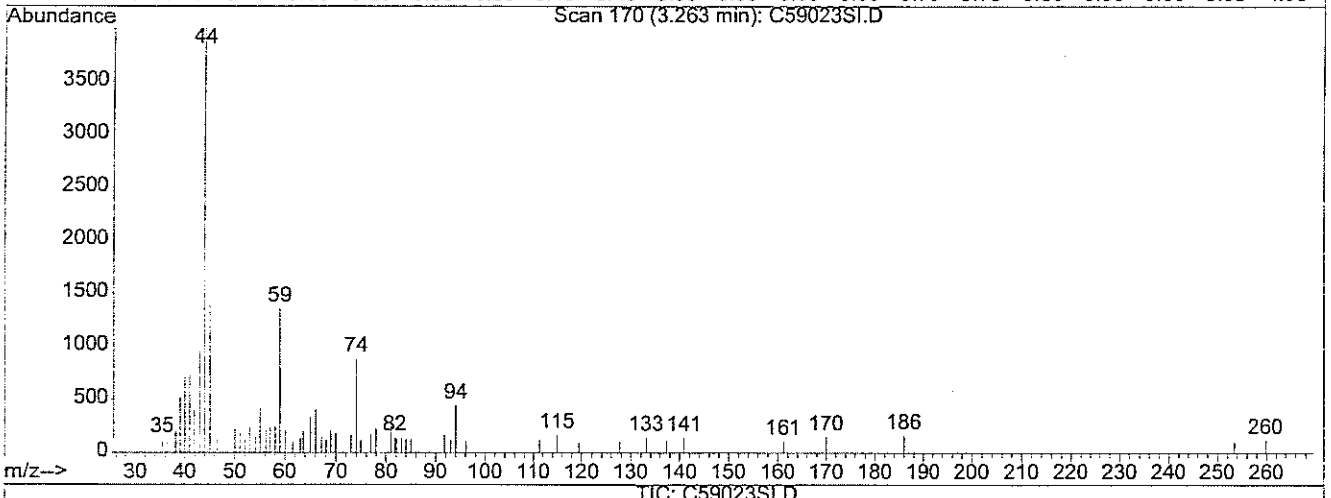
Quantitation Report (Qedit)

Data File : D:\HPCHEM\1\DATA\052206C\C59023SI.D Vial: 3
Acq On : 22 May 2006 11:24 am Operator:
Sample : 1 PPB Inst : Instr_C
Misc : 5000 Multiplr: 1.00
MS Integration Params: rteint.p
Quant Time: May 22 14:08 2006 Quant Results File: temp.res

Method : D:\HPCHEM\1\METHODS\V805226C.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Mon May 22 14:05:38 2006
Response via : Multiple Level Calibration

MT 5/23/06
extra area





TIC: C59023SI.D

(9) Diethyl ether (TM)

3.26min 1.42ug/L

response 3490

Ion	Exp%	Act%
74.00	100	100
45.00	149.50	92.38#
59.00	200.90	166.91
0.00	0.00	0.00

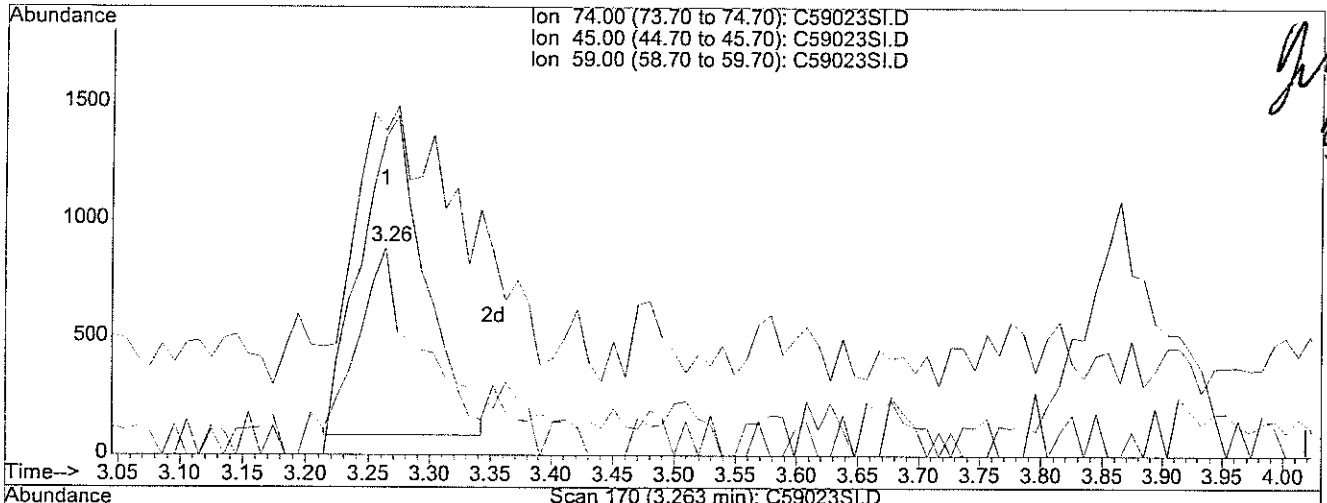
Quantitation Report (Qedit)

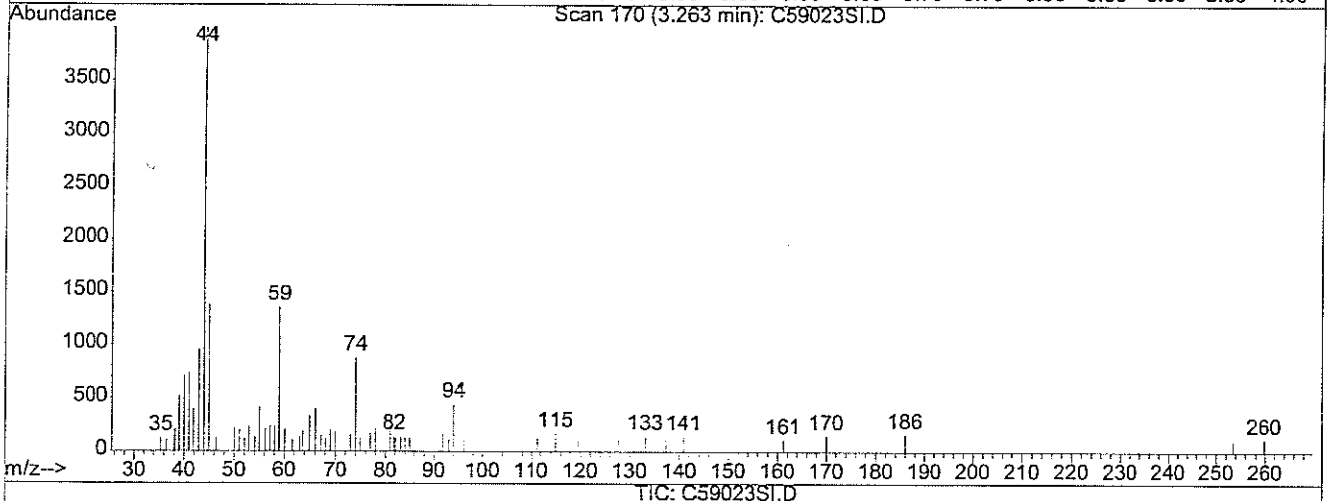
Data File : D:\HPCHEM\1\DATA\052206C\C59023SI.D
Acq On : 22 May 2006 11:24 am
Sample : 1 PPB
Misc : 5000
MS Integration Params: rteint.p
Quant Time: May 22 14:08 2006

Vial: 3
Operator:
Inst : Instr_C
Multiplr: 1.00
Quant Results File: temp.res

Method : D:\HPCHEM\1\METHODS\V805226C.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Mon May 22 14:05:38 2006
Response via : Multiple Level Calibration

MT 5/23/06





TIC: C59023SI.D

(9) Diethyl ether (TM)

3.26min 1.08ug/L m

response 2656

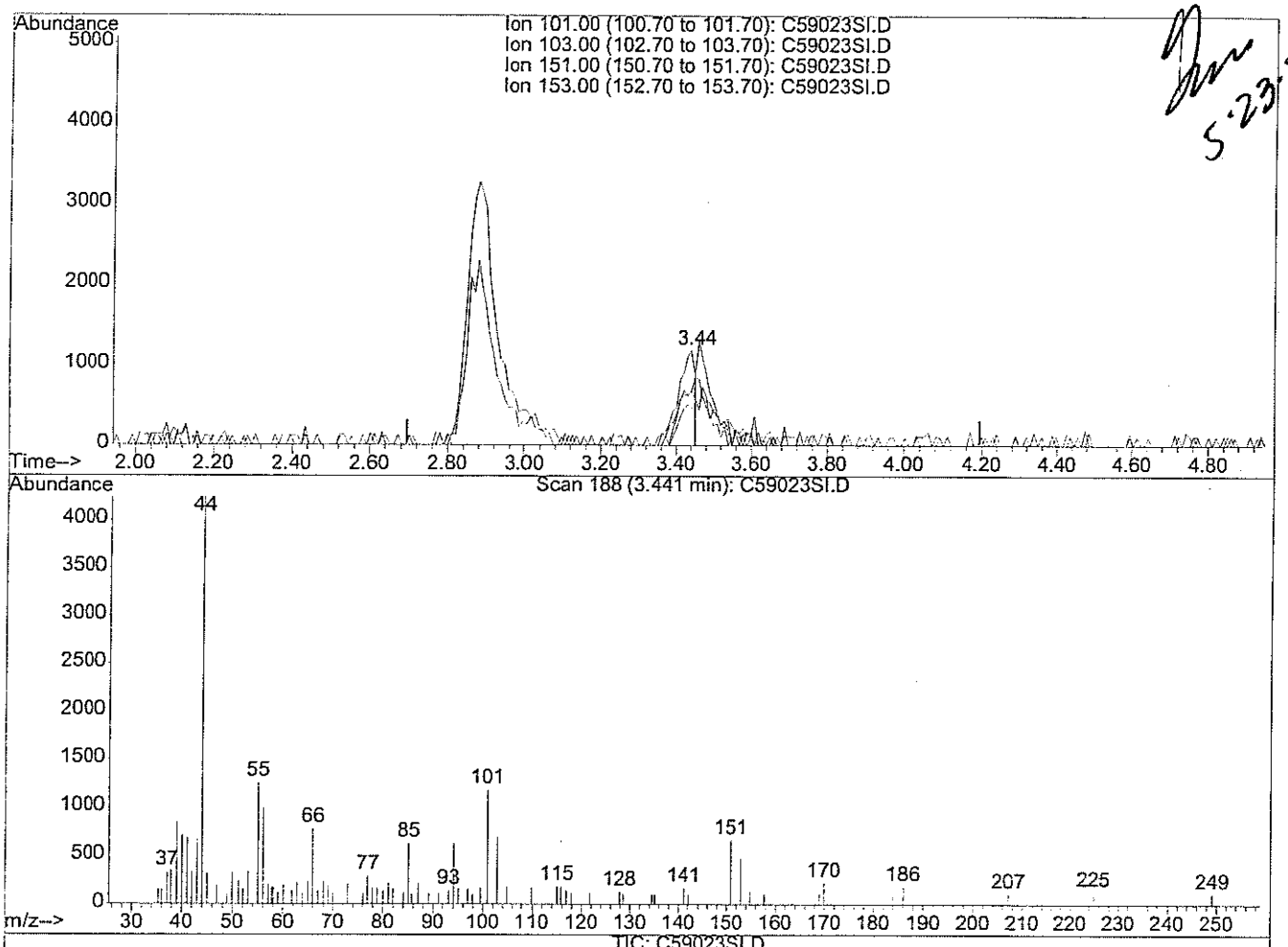
Ion	Exp%	Act%
74.00	100	100
45.00	149.50	121.39
59.00	200.90	219.31
0.00	0.00	0.00

Quantitation Report (Qedit)

Data File : D:\HPCHEM\1\DATA\052206C\C59023SI.D Vial: 3
Acq On : 22 May 2006 11:24 am Operator:
Sample : 1 PPB Inst : Instr_C
Misc : 5000 Multiplr: 1.00
MS Integration Params: rteint.p
Quant Time: May 22 13:18 2006 Quant Results File: temp.res

Method : D:\HPCHEM\1\METHODS\V805226C.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Wed Apr 19 15:41:22 2006
Response via : Multiple Level Calibration

MT 5/23/06
split peak



5.23.06

(10) 1,1,2-Trichlorotrifluoroethane (T)

3.44min 0.63ug/L

response 3868

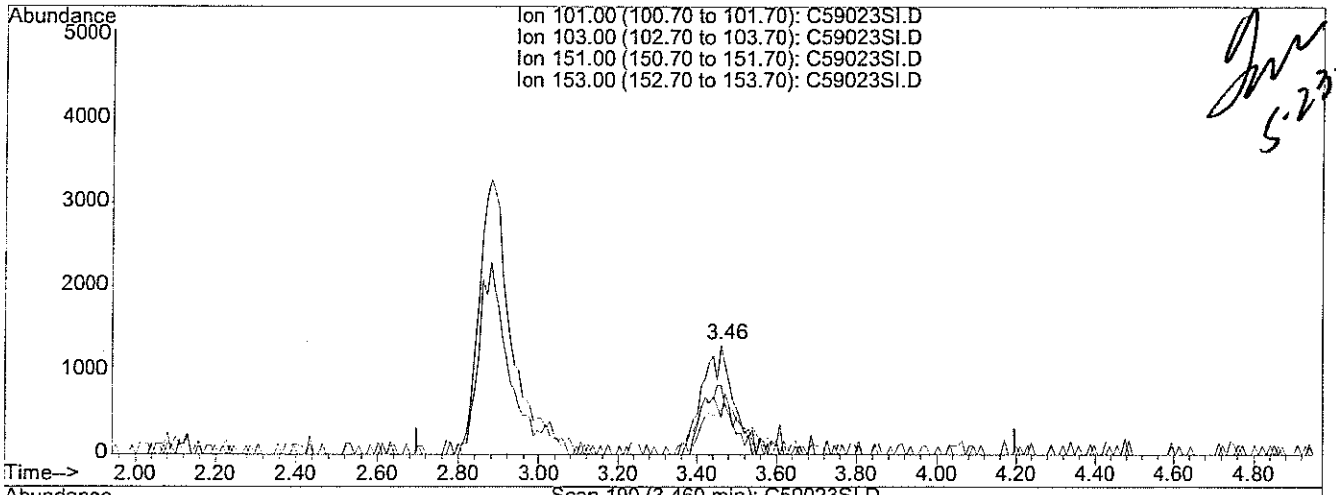
Ion	Exp%	Act%
101.00	100	100
103.00	65.50	73.32
151.00	74.70	124.53#
153.00	48.10	84.59#

Quantitation Report (Qedit)

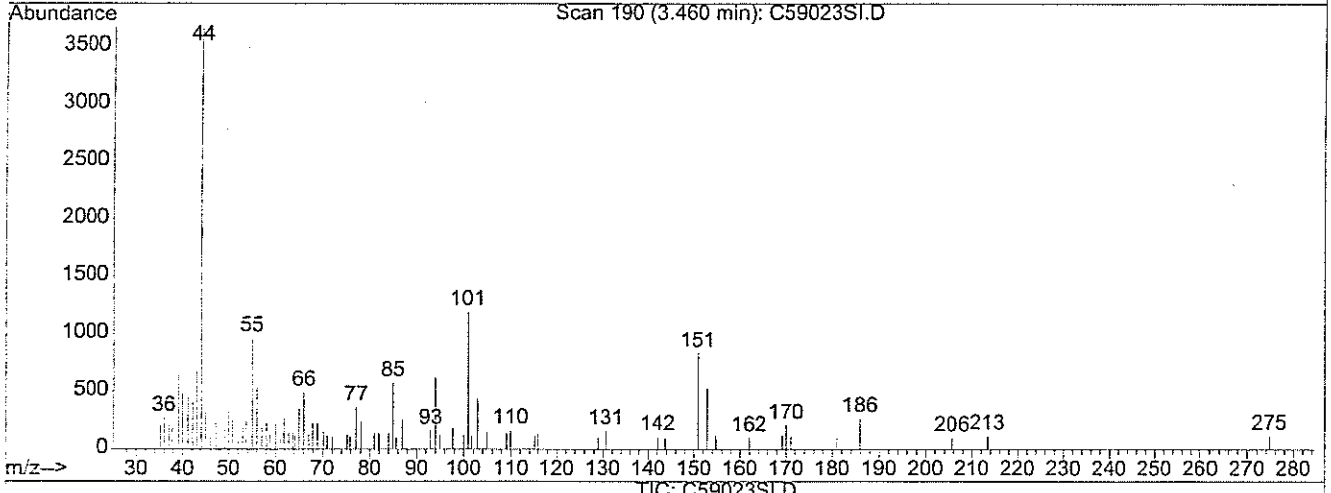
Data File : D:\HPCHEM\1\DATA\052206C\C59023SI.D Vial: 3
Acq On : 22 May 2006 11:24 am Operator:
Sample : 1 PPB Inst : Instr_C
Misc : 5000 Multiplr: 1.00
MS Integration Params: rteint.p
Quant Time: May 22 13:19 2006 Quant Results File: temp.res

Method : D:\HPCHEM\1\METHODS\V805226C.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Wed Apr 19 15:41:22 2006
Response via : Multiple Level Calibration

MT 5/23/06



5.23.06



TIC: C59023SI.D

(10) 1,1,2-Trichlorotrifluoroethane (T)

3.46min 1.25ug/L m

response 7665

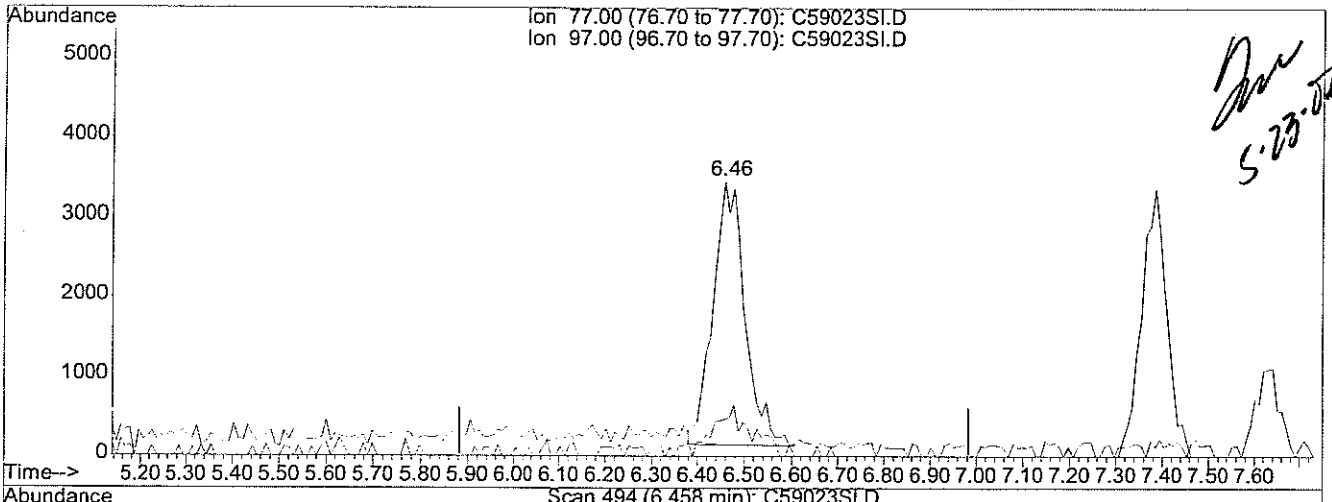
Ion	Exp%	Act%
101.00	100	100
103.00	65.50	37.00#
151.00	74.70	62.84
153.00	48.10	42.69

Quantitation Report (Qedit)

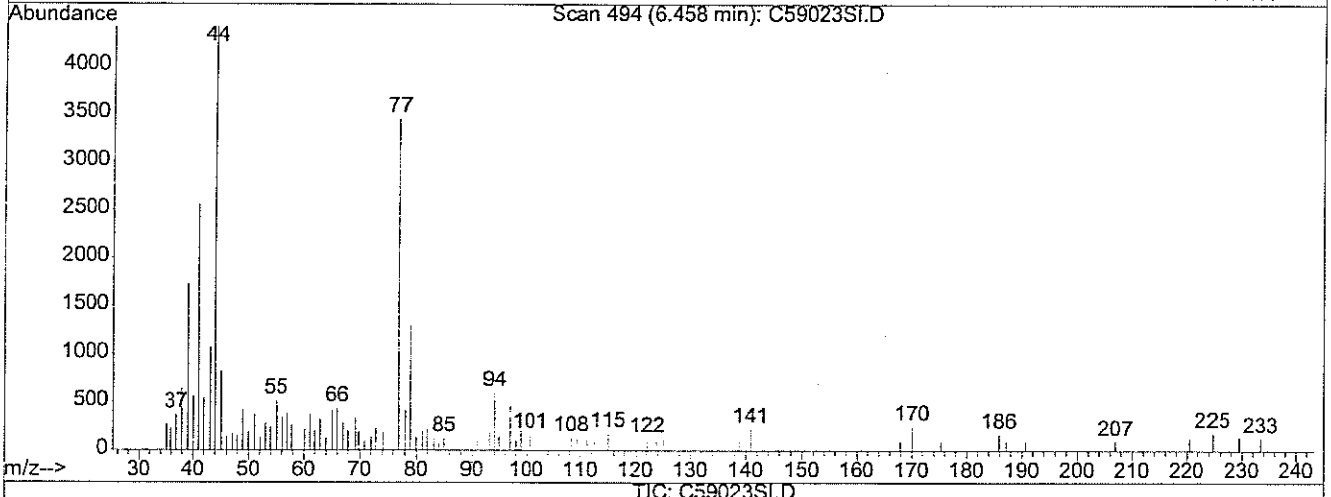
Data File : D:\HPCHEM\1\DATA\052206C\C59023SI.D Vial: 3
Acq On : 22 May 2006 11:24 am Operator:
Sample : 1 PPB Inst : Instr_C
Misc : 5000 Multiplr: 1.00
MS Integration Params: rteint.p
Quant Time: May 22 14:08 2006 Quant Results File: temp.res

Method : D:\HPCHEM\1\METHODS\V805226C.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Mon May 22 14:09:16 2006
Response via : Multiple Level Calibration

MT 5/23/06
extra area



5.23.06



(25) 2,2-Dichloropropane (TM)

6.46min 1.63ug/L

response 15670

Ion	Exp%	Act%
77.00	100	100
97.00	19.40	14.00
0.00	0.00	0.00
0.00	0.00	0.00

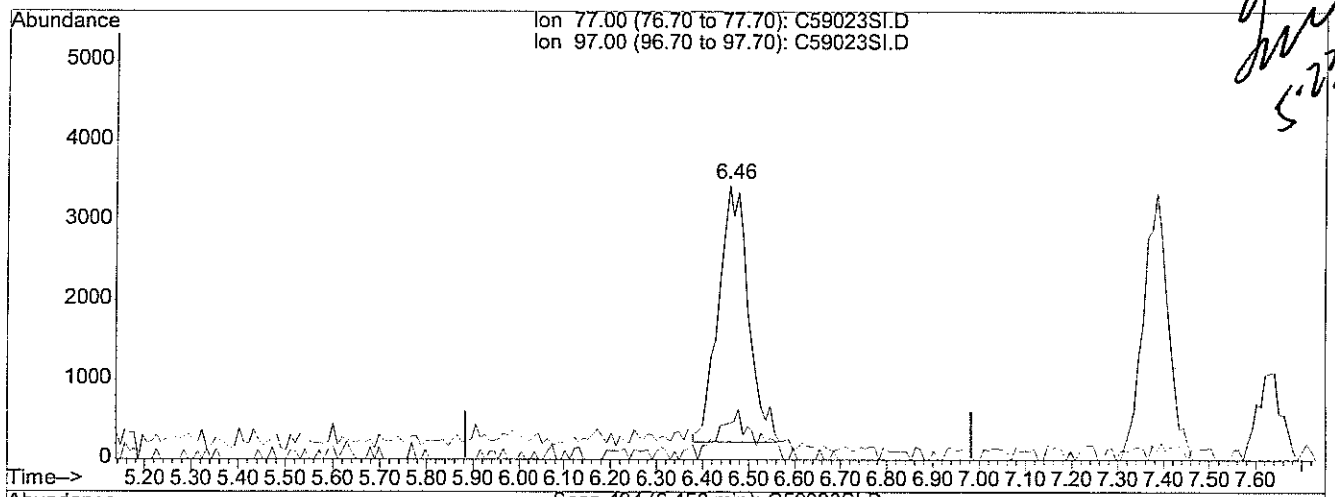
Quantitation Report (Qedit)

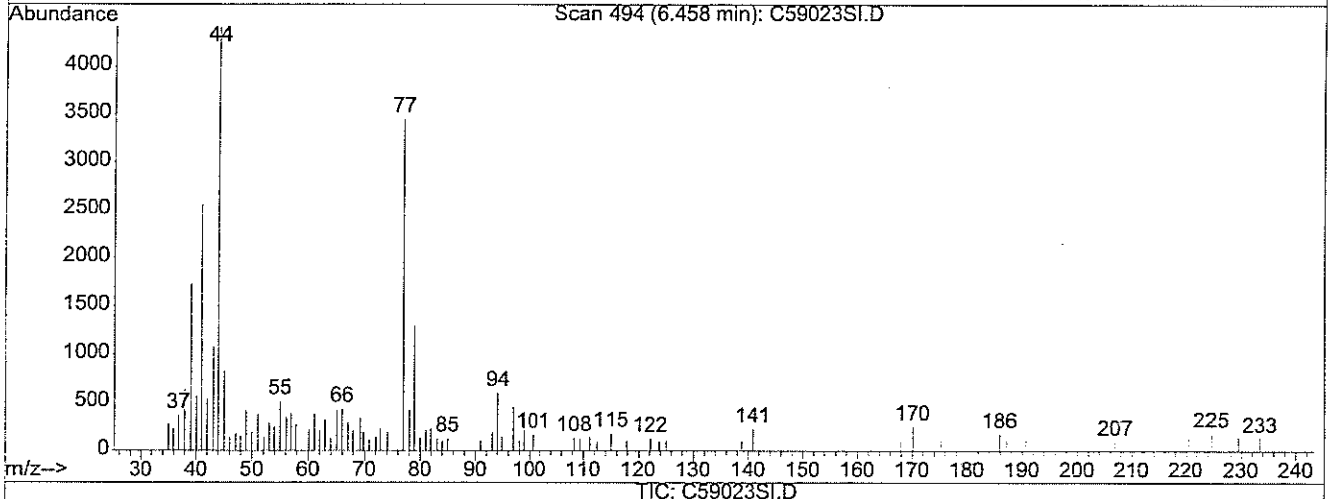
Data File : D:\HPCHEM\1\DATA\052206C\C59023SI.D Vial: 3
Acq On : 22 May 2006 11:24 am Operator:
Sample : 1 PPB Inst : Instr_C
Misc : 5000 Multiplr: 1.00
MS Integration Params: rteint.p
Quant Time: May 22 14:11 2006 Quant Results File: temp.res

Method : D:\HPCHEM\1\METHODS\V805226C.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Mon May 22 14:18:09 2006
Response via : Multiple Level Calibration

MT 5/23/06

5.23.06





(25) 2,2-Dichloropropane (TM)

6.46min 1.52ug/L m

response 14594

Ion	Exp%	Act%
77.00	100	100
97.00	19.40	13.38
0.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (QT Reviewed)

Data File : D:\HPCHEM\1\DATA\052206C\C59024SI.D Vial: 4
Acq On : 22 May 2006 11:57 am Operator:
Sample : 2 PPB Inst : Instr_C
Misc : 5000 Multiplr: 1.00
Integration Params: rteint.p
Quant Time: May 22 14:08 2006 Quant Results File: V805226C.RES

Quant Method : D:\HPCHEM\1\METHODS\V805226C.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Wed Apr 19 15:41:22 2006
Response via : Initial Calibration
DataAcq Meth : V805226C

MT 5/23/06 5.23.06

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.00	168	257975	20.00	ug/L	-0.01
39) 1,4-Difluorobenzene	8.48	114	426815 ✓	20.00	ug/L	-0.01
63) *d5-Chlorobenzene	13.02	117	375633 ✓	20.00	ug/L	0.00
87) 1,4-Dichlorobenzene-d4	16.74	152	188191	20.00	ug/L	0.00
System Monitoring Compounds						
28) DB Holder #28	7.86	65	18591	1.81	ug/L	-0.01
Spiked Amount 20.000	Range 70 - 130		Recovery = 9.05%#			
36) d4-1,2-Dichloroethane	7.86	65	18591	1.81	ug/L	-0.01
Spiked Amount 20.000	Range 70 - 130		Recovery = 9.05%#			
54) d8-Toluene	10.74	98	42292	1.77	ug/L	0.00
Spiked Amount 20.000	Range 70 - 130		Recovery = 8.85%#			
73) Bromofluorobenzene	14.89	95	18280	1.92	ug/L	0.00
Spiked Amount 20.000	Range 70 - 130		Recovery = 9.60%#			
Target Compounds						Qvalue
2) Dichlorodifluoromethane	1.69	85	13723	1.58	ug/L	87
3) Chloromethane	1.94	50	13692	1.57	ug/L	95
4) Vinyl Chloride	2.02	62	15473	1.73	ug/L	85
5) Bromomethane	2.48	94	6135	1.33	ug/L	95
6) Chloroethane	2.57	64	8220	1.88	ug/L	99
7) t-Butyl alcohol (TBA)	3.88	59	8841	13.20	ug/L #	82
8) Trichlorofluoromethane	2.87	101	26905	1.67	ug/L	96
9) Diethyl ether	3.27	74	6220	2.45	ug/L	87
10) 1,1,2-Trichlorotrifluoroet	3.45	101	14760	2.34	ug/L	96
11) Acrolein	3.41	56	10739	19.27	ug/L #	76
12) Acetone	3.52	43	25143m	11.23	ug/L	
13) 1,1-Dichloroethene	3.67	96	10929	2.38	ug/L	87
14) Methyl iodide	4.10	142	17213	2.42	ug/L	92
15) Di-isopropyl ether (DIPE)	5.49	45	48468	2.14	ug/L	96
16) Methylene Chloride	4.41	84	13551	2.71	ug/L #	87
17) Carbon Disulfide	4.38	76	35186	2.39	ug/L	100
18) Acrylonitrile	4.66	53	4846	2.29	ug/L #	88
19) Methyl-tert-butyl ether (M	4.73	73	41701	2.41	ug/L	100
20) trans-1,2-Dichloroethene	4.89	61	25337	2.25	ug/L	96
21) 1,1-Dichloroethane	5.58	63	29177	2.27	ug/L	91
22) Vinyl acetate	5.63	43	32464	3.02	ug/L	99
23) Methyl ethyl ketone	6.30	72	5452	9.96	ug/L #	27
24) Ethyl t-butyl ether (ETBE)	6.14	59	49921	2.35	ug/L	98
25) 2,2-Dichloropropane	6.47	77	21506	2.17	ug/L	93
26) cis-1,2-Dichloroethene	6.53	96	16576	2.27	ug/L	87
27) t-Amyl methyl ether (TAME)	7.87	73	35587	2.40	ug/L	93

(#) = qualifier out of range (m) = manual integration

Quantitation Report (QT Reviewed)

Data File : D:\HPCHEM\1\DATA\052206C\C59024SI.D Vial: 4
Acq On : 22 May 2006 11:57 am Operator:
Sample : 2 PPB Inst : Instr_C
Misc : 5000 Multiplr: 1.00
Integration Params: rteint.p
Quant Time: May 22 14:08 2006 Quant Results File: V805226C.RES

Quant Method : D:\HPCHEM\1\METHODS\V805226C.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Wed Apr 19 15:41:22 2006
Response via : Initial Calibration
DataAcq Meth : V805226C

	Compound	R.T.	QIon	Response	Conc	Unit		Qvalue
29)	Chloroform	6.78	83	27246	2.24	ug/L		93
30)	Bromochloromethane	7.01	128	9278	2.39	ug/L		94
31)	Tetrahydrofuran	7.13	71	1446	2.44	ug/L	#	60
32)	1,1,1-Trichloroethane	7.38	97	23618	2.14	ug/L		88
34)	1,1-Dichloropropene	7.62	75	20551	2.34	ug/L		98
35)	Carbon Tetrachloride	7.75	119	19165	2.10	ug/L		96
37)	1,2-Dichloroethane	8.01	62	25823	2.16	ug/L		96
38)	Benzene	8.01	78	49606	2.37	ug/L		93
40)	Trichloroethene	8.95	95	16940	2.26	ug/L		99
41)	1,2-Dichloropropane	9.21	63	16458	2.33	ug/L		99
42)	Methylmethacrylate	9.37	69	9378	2.14	ug/L		92
43)	Methyl acetate	4.22	43	12478	1.93	ug/L	#	93
44)	Cyclohexane	7.39	56	26390	2.31	ug/L		96
45)	Bromodichloromethane	9.56	83	18524	2.01	ug/L		94
46)	Dibromomethane	9.63	93	11618	2.15	ug/L		92
47)	1,4-Dioxane	9.64	88	4613	69.10	ug/L		91
48)	Methylcyclohexane	9.03	83	21823	2.39	ug/L		98
50)	2-Chloroethylvinyl ether	10.08	63	3657	1.75	ug/L		91
51)	2-Hexanone	11.44	58	29855	12.69	ug/L		92
52)	Methyl isobutyl ketone	10.13	58	30427	12.54	ug/L		96
53)	cis-1,3-Dichloropropene	10.39	75	22341	2.13	ug/L		89
55)	Toluene	10.86	92	33130	2.34	ug/L		99
56)	trans-1,3-Dichloropropene	11.15	75	21047	2.10	ug/L		93
58)	1,1,2-Trichloroethane	11.37	83	11309	2.43	ug/L		92
59)	1,3-Dichloropropane	11.76	76	21124	2.32	ug/L		98
60)	Tetrachloroethene	11.84	166	19947	2.19	ug/L		98
61)	Dibromochloromethane	12.12	129	15404	1.96	ug/L		93
62)	1,2-Dibromoethane	12.43	107	18170	2.31	ug/L		97
64)	Chlorobenzene	13.07	112	43066	2.27	ug/L		95
65)	1,1,1,2-Tetrachloroethane	13.15	131	15167	2.26	ug/L		92
66)	Ethylbenzene	13.18	91	61570	2.25	ug/L		98
67)	m,p-Xylene	13.29	106	52018	4.74	ug/L		97
68)	o-Xylene	13.96	106	25372	2.39	ug/L		93
69)	Styrene	14.01	104	43617	2.34	ug/L		92
70)	Bromoform	14.49	173	11764	2.02	ug/L		89
71)	Isopropylbenzene	14.52	105	64923	2.30	ug/L		99
72)	1,1,2,2-Tetrachloroethane	14.78	83	17178	2.33	ug/L		99
74)	1,2,3-Trichloropropane	15.00	75	12078	2.21	ug/L		91
75)	trans-1,4-Dichloro-2-buten	15.14	89	5059	2.22	ug/L		98
76)	n-Propylbenzene	15.14	91	73695	2.34	ug/L		98
77)	Bromobenzene	15.17	156	18197	2.24	ug/L		97

(#) = qualifier out of range (m) = manual integration

Quantitation Report (QT Reviewed)

Data File : D:\HPCHEM\1\DATA\052206C\C59024SI.D Vial: 4
Acq On : 22 May 2006 11:57 am Operator:
Sample : 2 PPB Inst : Instr_C
Misc : 5000 Multiplr: 1.00
Integration Params: rteint.p
Quant Time: May 22 14:08 2006 Quant Results File: V805226C.RES

Quant Method : D:\HPCHEM\1\METHODS\V805226C.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Wed Apr 19 15:41:22 2006
Response via : Initial Calibration
DataAcq Meth : V805226C

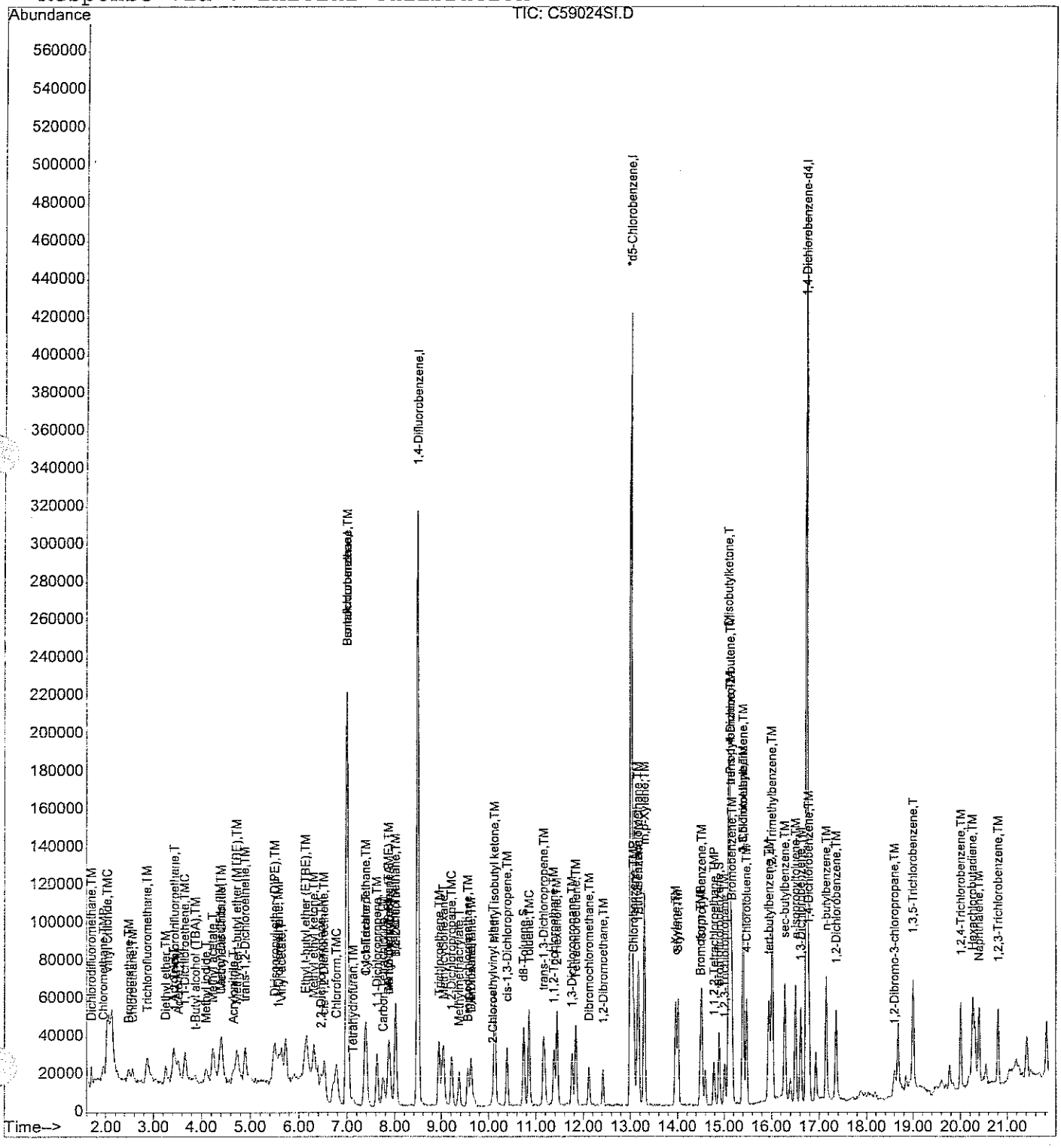
	Compound	R.T.	QIon	Response	Conc	Unit		Qvalue
78)	1,3,5-Trimethylbenzene	15.39	105	51890	2.23	ug/L		99
79)	2-Chlorotoluene	15.39	91	46368	2.13	ug/L		98
80)	4-Chlorotoluene	15.47	91	45958	2.30	ug/L		94
81)	tert-butylbenzene	15.95	119	45964	2.24	ug/L		95
82)	1,2,4-Trimethylbenzene	16.01	105	55908	2.36	ug/L		98
83)	Diisobutylketone	15.10	57	188240	11.37	ug/L		91
84)	sec-butylbenzene	16.28	105	72264	2.28	ug/L		94
85)	p-isopropyltoluene	16.51	119	57834	2.27	ug/L		96
86)	1,3-Dichlorobenzene	16.62	146	30219	2.20	ug/L		94
88)	1,4-Dichlorobenzene	16.79	146	31299	2.23	ug/L		97
89)	n-butylbenzene	17.15	91	52067	2.21	ug/L		96
90)	1,2-Dichlorobenzene	17.36	146	27744	2.14	ug/L		99
92)	1,2-Dibromo-3-chloropropan	18.61	75	3161	2.28	ug/L	#	66
94)	1,2,4-Trichlorobenzene	20.01	180	22584	2.03	ug/L		95
95)	Hexachlorobutadiene	20.26	225	11196	1.97	ug/L		95
96)	Naphthalene	20.39	128	43368	2.13	ug/L		100
97)	1,2,3-Trichlorobenzene	20.81	180	20074	1.99	ug/L		89
98)	1,3,5-Trichlorobenzene	18.99	180	26039	2.08	ug/L		86

(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : D:\HPCHEM\1\DATA\052206C\C59024SI.D Vial: 4
Acq On : 22 May 2006 11:57 am Operator:
Sample : 2 PPB Inst : Instr_C
Misc : 5000 Multiplr: 1.00
MS Integration Params: rteint.p
Quant Time: May 22 14:08 2006 Quant Results File: V805226C.RES

Method : D:\HPCHEM\1\METHODS\V805226C.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Mon May 22 14:18:09 2006
Response via : Initial Calibration



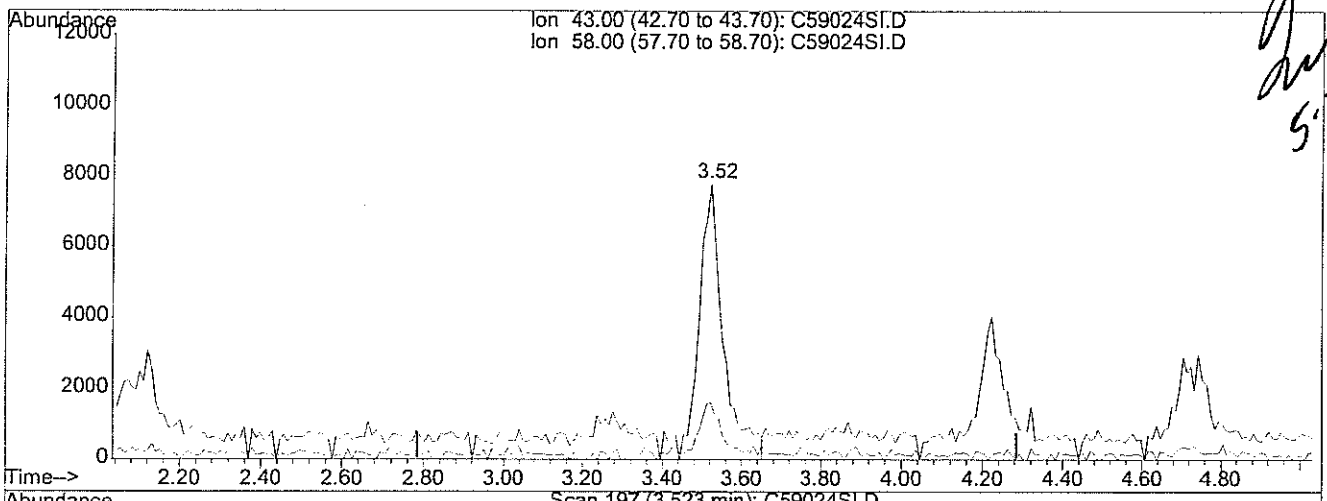
Quantitation Report (Qedit)

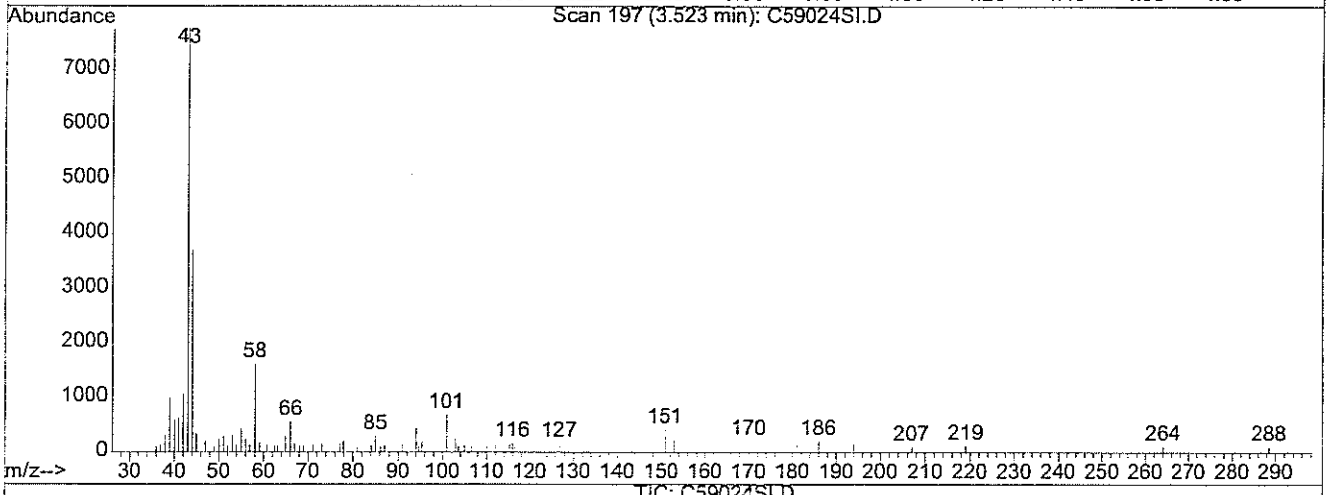
Data File : D:\HPCHEM\1\DATA\052206C\C59024SI.D Vial: 4
Acq On : 22 May 2006 11:57 am Operator:
Sample : 2 PPB Inst : Instr_C
Misc : 5000 Multiplr: 1.00
MS Integration Params: rteint.p
Quant Time: May 22 13:18 2006 Quant Results File: temp.res

Method : D:\HPCHEM\1\METHODS\V805226C.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Mon May 22 14:05:38 2006
Response via : Multiple Level Calibration

MT 5/23/06
extra area

5.23.06





TIC: C59024SI.D

(12) Acetone (TM)

3.52min 14.78ug/L

response 33083

Ion	Exp%	Act%
43.00	100	100
58.00	19.60	16.47
0.00	0.00	0.00
0.00	0.00	0.00

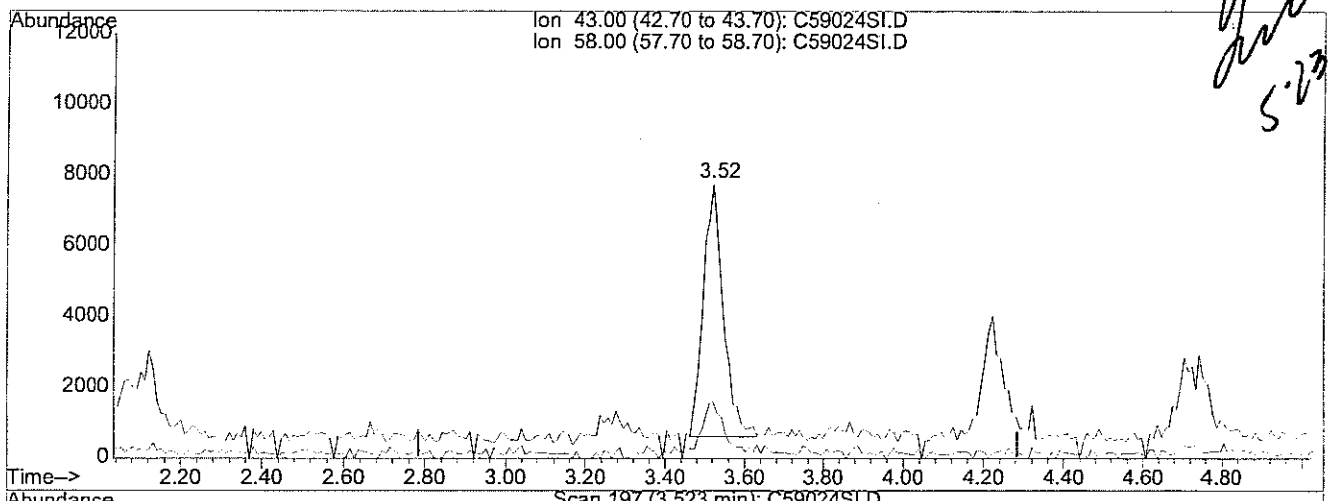
Quantitation Report (Qedit)

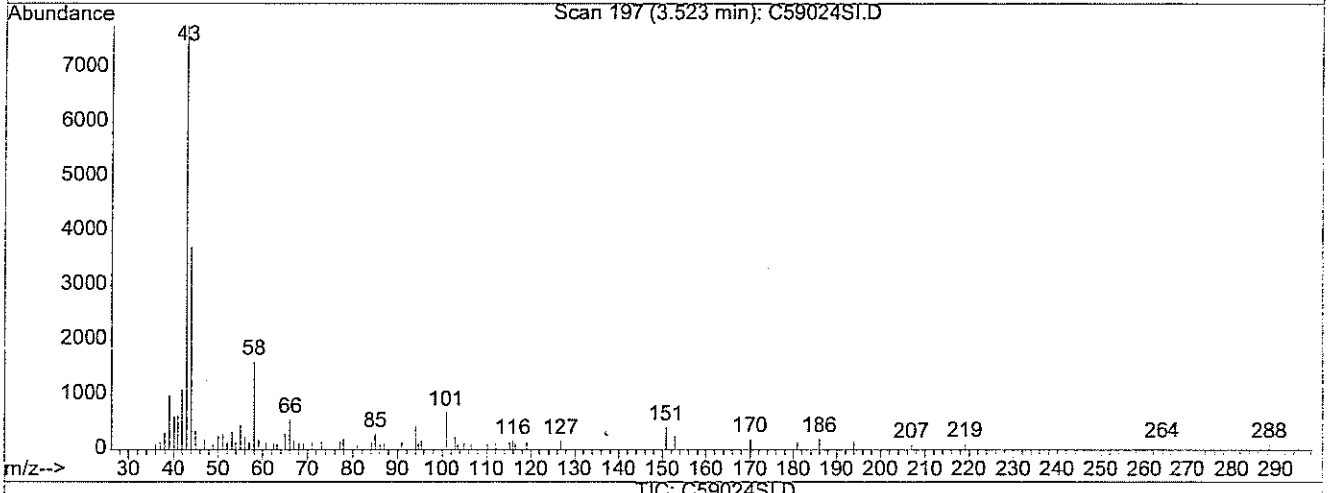
Data File : D:\HPCHEM\1\DATA\052206C\C59024SI.D Vial: 4
Acq On : 22 May 2006 11:57 am Operator:
Sample : 2 PPB Inst : Instr_C
Misc : 5000 Multiplr: 1.00
MS Integration Params: rteint.p
Quant Time: May 22 14:08 2006 Quant Results File: temp.res

Method : D:\HPCHEM\1\METHODS\V805226C.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Mon May 22 14:05:38 2006
Response via : Multiple Level Calibration

MT 5/23/06

5.23.06





TIC: C59024SI.D

(12) Acetone (TM)

3.52min 11.23ug/L m

response 25143

Ion	Exp%	Act%
43.00	100	100
58.00	19.60	21.67
0.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (QT Reviewed)

Data File : D:\HPCHEM\1\DATA\052206C\C59025SI.D Vial: 5
Acq On : 22 May 2006 12:30 pm Operator:
Sample : 5 PPB Inst : Instr_C
Misc : 5000 Multiplr: 1.00
Integration Params: rteint.p
Quant Time: May 22 13:18 2006 Quant Results File: V805226C.RES

Quant Method : D:\HPCHEM\1\METHODS\V805226C.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Wed Apr 19 15:41:22 2006
Response via : Initial Calibration
DataAcq Meth : V805226C

MT 5/23/06 5.27.06

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.00	168	251193	20.00	ug/L	-0.02
39) 1,4-Difluorobenzene	8.48	114	421056 ✓	20.00	ug/L	-0.01
63) *d5-Chlorobenzene	13.01	117	376549 ✓	20.00	ug/L	-0.01
87) 1,4-Dichlorobenzene-d4	16.74	152	181024	20.00	ug/L	0.00
System Monitoring Compounds						
28) DB Holder #28	7.87	65	45609	4.57	ug/L	0.00
Spiked Amount 20.000	Range 70 - 130		Recovery =	22.85%#		
36) d4-1,2-Dichloroethane	7.87	65	45609	4.57	ug/L	0.00
Spiked Amount 20.000	Range 70 - 130		Recovery =	22.85%#		
54) d8-Toluene	10.74	98	111927	4.75	ug/L	0.00
Spiked Amount 20.000	Range 70 - 130		Recovery =	23.75%#		
73) Bromofluorobenzene	14.89	95	44299	4.63	ug/L	0.00
Spiked Amount 20.000	Range 70 - 130		Recovery =	23.15%#		
Target Compounds						Qvalue
2) Dichlorodifluoromethane	1.69	85	36361	4.29	ug/L	94
3) Chloromethane	1.96	50	35888	4.22	ug/L	100
4) Vinyl Chloride	2.03	62	40767	4.69	ug/L	98
5) Bromomethane	2.48	94	23004	5.11	ug/L	95
6) Chloroethane	2.57	64	20344	4.77	ug/L	95
7) t-Butyl alcohol (TBA)	3.87	59	18903	28.99	ug/L #	77
8) Trichlorofluoromethane	2.87	101	73064	4.65	ug/L	99
9) Diethyl ether	3.26	74	12961	5.24	ug/L	88
10) 1,1,2-Trichlorotrifluoroet	3.45	101	32000	5.22	ug/L	99
11) Acrolein	3.41	56	20529	37.84	ug/L	97
12) Acetone	3.52	43	59030	27.08	ug/L	96
13) 1,1-Dichloroethene	3.66	96	24460	5.48	ug/L	95
14) Methyl iodide	4.11	142	41140	5.93	ug/L	96
15) Di-isopropyl ether (DIPE)	5.50	45	107043	4.86	ug/L	99
16) Methylene Chloride	4.41	84	26828	5.51	ug/L	93
17) Carbon Disulfide	4.37	76	78493	5.48	ug/L	100
18) Acrylonitrile	4.64	53	12394	6.00	ug/L	90
19) Methyl-tert-butyl ether (M	4.73	73	89006	5.27	ug/L	98
20) trans-1,2-Dichloroethene	4.89	61	56401	5.14	ug/L	99
21) 1,1-Dichloroethane	5.59	63	64195	5.12	ug/L	99
22) Vinyl acetate	5.63	43	62179	5.95	ug/L	96
23) Methyl ethyl ketone	6.31	72	15511	29.10	ug/L	84
24) Ethyl t-butyl ether (ETBE)	6.15	59	111241	5.39	ug/L	98
25) 2,2-Dichloropropane	6.45	77	52062	5.41	ug/L	89
26) cis-1,2-Dichloroethene	6.53	96	38177	5.36	ug/L	92
27) t-Amyl methyl ether (TAME)	7.87	73	77745	5.38	ug/L	96

(#) = qualifier out of range (m) = manual integration

Quantitation Report (QT Reviewed)

Data File : D:\HPCHEM\1\DATA\052206C\C59025SI.D Vial: 5
Acq On : 22 May 2006 12:30 pm Operator:
Sample : 5 PPB Inst : Instr_C
Misc : 5000 Multiplr: 1.00
Integration Params: rteint.p
Quant Time: May 22 13:18 2006 Quant Results File: V805226C.RES

Quant Method : D:\HPCHEM\1\METHODS\V805226C.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Wed Apr 19 15:41:22 2006
Response via : Initial Calibration
DataAcq Meth : V805226C

	Compound	R.T.	QIon	Response	Conc	Unit		Qvalue
29)	Chloroform	6.78	83	60244	5.10	ug/L		94
30)	Bromochloromethane	7.01	128	19918	5.28	ug/L		88
31)	Tetrahydrofuran	7.12	71	3721	6.46	ug/L		95
32)	1,1,1-Trichloroethane	7.38	97	54496	5.07	ug/L		97
34)	1,1-Dichloropropene	7.63	75	44970	5.25	ug/L		95
35)	Carbon Tetrachloride	7.75	119	45107	5.08	ug/L		98
37)	1,2-Dichloroethane	8.01	62	57407	4.92	ug/L		99
38)	Benzene	8.01	78	109686	5.38	ug/L		99
40)	Trichloroethene	8.95	95	38027	5.14	ug/L		95
41)	1,2-Dichloropropane	9.21	63	35988	5.16	ug/L		96
42)	Methylmethacrylate	9.36	69	26059	6.04	ug/L		93
43)	Methyl acetate	4.22	43	33979	5.34	ug/L	#	96
44)	Cyclohexane	7.38	56	58113	5.16	ug/L		98
45)	Bromodichloromethane	9.56	83	43523	4.78	ug/L		96
46)	Dibromomethane	9.63	93	28051	5.27	ug/L		92
47)	1,4-Dioxane	9.64	88	10269	155.93	ug/L		96
48)	Methylcyclohexane	9.04	83	48534	5.40	ug/L		93
50)	2-Chloroethylvinyl ether	10.07	63	7909	3.83	ug/L	#	88
51)	2-Hexanone	11.43	58	66621	28.70	ug/L		95
52)	Methyl isobutyl ketone	10.13	58	68974	28.82	ug/L		99
53)	cis-1,3-Dichloropropene	10.40	75	50773	4.90	ug/L		96
55)	Toluene	10.86	92	72241	5.17	ug/L		94
56)	trans-1,3-Dichloropropene	11.16	75	48335	4.90	ug/L		97
58)	1,1,2-Trichloroethane	11.37	83	24390	5.31	ug/L		91
59)	1,3-Dichloropropane	11.76	76	47603	5.30	ug/L		95
60)	Tetrachloroethene	11.84	166	43432	4.84	ug/L		99
61)	Dibromochloromethane	12.12	129	35074	4.53	ug/L		99
62)	1,2-Dibromoethane	12.42	107	40968	5.27	ug/L		95
64)	Chlorobenzene	13.07	112	95455	5.03	ug/L		98
65)	1,1,1,2-Tetrachloroethane	13.15	131	32118	4.77	ug/L		97
66)	Ethylbenzene	13.17	91	137871	5.03	ug/L		100
67)	m,p-Xylene	13.30	106	114020	10.37	ug/L		92
68)	o-Xylene	13.96	106	53292	5.01	ug/L		98
69)	Styrene	14.01	104	95246	5.09	ug/L		99
70)	Bromoform	14.49	173	28335	4.85	ug/L		94
71)	Isopropylbenzene	14.52	105	142773	5.05	ug/L		99
72)	1,1,2,2-Tetrachloroethane	14.79	83	39009	5.28	ug/L		92
74)	1,2,3-Trichloropropane	15.00	75	29641	5.40	ug/L	#	90
75)	trans-1,4-Dichloro-2-buten	15.13	89	11727	5.13	ug/L		96
76)	n-Propylbenzene	15.14	91	161911	5.12	ug/L		98
77)	Bromobenzene	15.17	156	40927	5.03	ug/L		97

(#) = qualifier out of range (m) = manual integration

Quantitation Report (QT Reviewed)

Data File : D:\HPCHEM\1\DATA\052206C\C59025SI.D Vial: 5
Acq On : 22 May 2006 12:30 pm Operator:
Sample : 5 PPB Inst : Instr_C
Misc : 5000 Multiplr: 1.00
Integration Params: rteint.p
Quant Time: May 22 13:18 2006 Quant Results File: V805226C.RES

Quant Method : D:\HPCHEM\1\METHODS\V805226C.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Wed Apr 19 15:41:22 2006
Response via : Initial Calibration
DataAcq Meth : V805226C

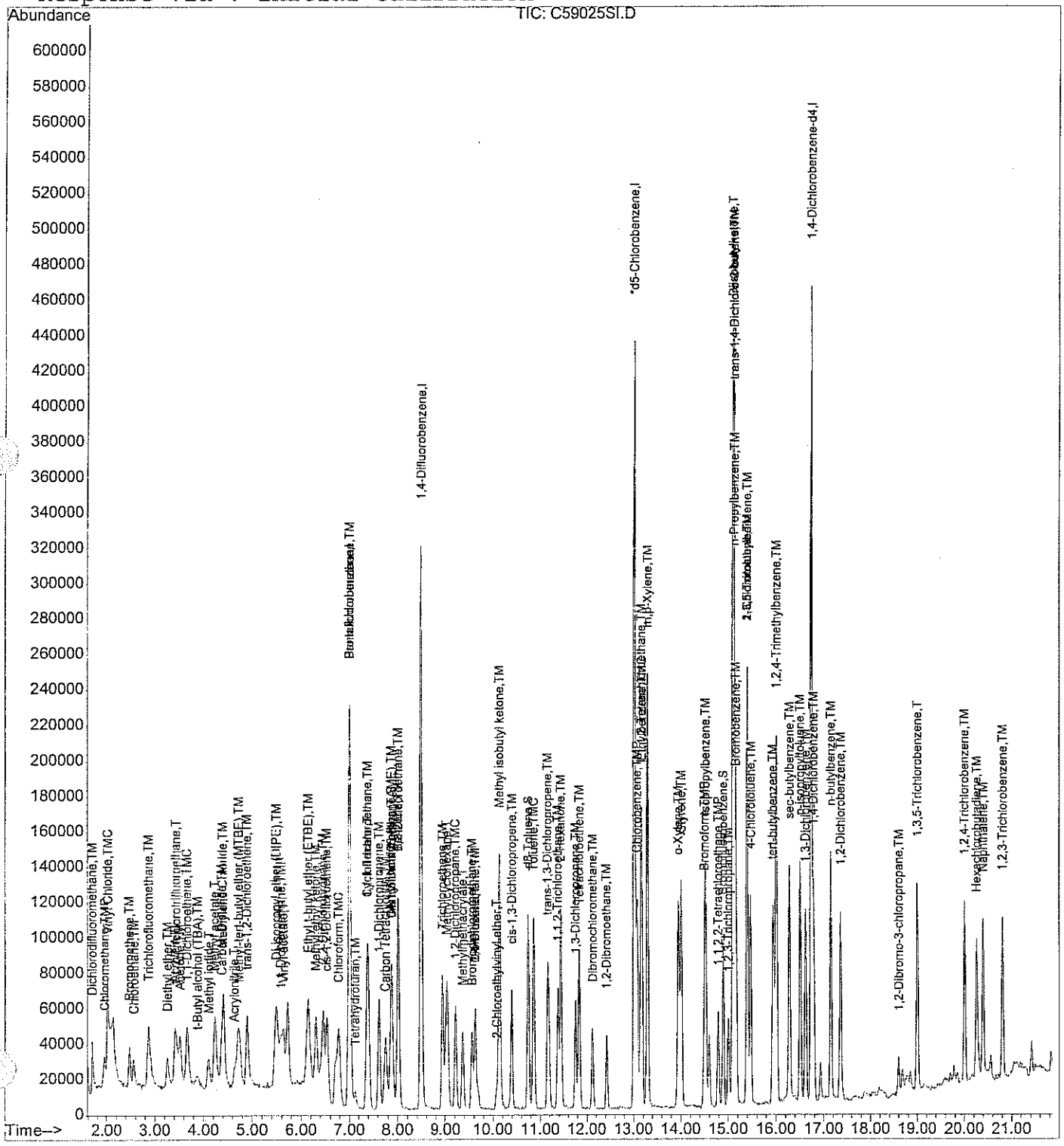
	Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
78)	1,3,5-Trimethylbenzene	15.40	105	116300	4.99	ug/L	93
79)	2-Chlorotoluene	15.40	91	107466	4.93	ug/L	99
80)	4-Chlorotoluene	15.48	91	96559	4.82	ug/L	100
81)	tert-butylbenzene	15.95	119	101082	4.91	ug/L	97
82)	1,2,4-Trimethylbenzene	16.01	105	119535	5.04	ug/L	95
83)	Diisobutylketone	15.10	57	411413	24.78	ug/L	93
84)	sec-butylbenzene	16.28	105	159303	5.02	ug/L	97
85)	p-isopropyltoluene	16.50	119	125437	4.92	ug/L	96
86)	1,3-Dichlorobenzene	16.62	146	67357	4.90	ug/L	98
88)	1,4-Dichlorobenzene	16.79	146	71448	5.28	ug/L	97
89)	n-butylbenzene	17.15	91	114401	5.05	ug/L	99
90)	1,2-Dichlorobenzene	17.36	146	63914	5.12	ug/L	98
92)	1,2-Dibromo-3-chloropropan	18.60	75	6815	5.11	ug/L	95
94)	1,2,4-Trichlorobenzene	20.01	180	51813	4.83	ug/L	97
95)	Hexachlorobutadiene	20.26	225	23170	4.23	ug/L	99
96)	Naphthalene	20.40	128	101093	5.17	ug/L	100
97)	1,2,3-Trichlorobenzene	20.81	180	47449	4.88	ug/L	94
98)	1,3,5-Trichlorobenzene	19.00	180	56838	4.72	ug/L	97

(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : D:\HPCHEM\1\DATA\052206C\C59025SI.D Vial: 5
Acq On : 22 May 2006 12:30 pm Operator:
Sample : 5 PPB Inst : Instr_C
Misc : 5000 Multiplr: 1.00
MS Integration Params: rteint.p
Quant Time: May 22 13:18 2006 Quant Results File: V805226C.RES

Method : D:\HPCHEM\1\METHODS\V805226C.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Mon May 22 14:18:09 2006
Response via : Initial Calibration



Quantitation Report (QT Reviewed)

Data File : D:\HPCHEM\1\DATA\052206C\C59026SI.D Vial: 6
Acq On : 22 May 2006 1:07 pm Operator:
Sample : 20 PPB Inst : Instr_C
Misc : 5000 Multiplr: 1.00
Integration Params: rteint.p
Quant Time: May 22 13:18 2006 Quant Results File: V805226C.RES

Quant Method : D:\HPCHEM\1\METHODS\V805226C.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Wed Apr 19 15:41:22 2006
Response via : Initial Calibration
DataAcq Meth : V805226C

MT 5/23/06 5.23.06

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.00	168	253263	20.00	ug/L	-0.02
39) 1,4-Difluorobenzene	8.48	114	428236 ✓	20.00	ug/L	-0.02
63) *d5-Chlorobenzene	13.02	117	378997 ✓	20.00	ug/L	0.00
87) 1,4-Dichlorobenzene-d4	16.73	152	180603	20.00	ug/L	0.00
System Monitoring Compounds						
28) DB Holder #28	7.87	65	187003	18.58	ug/L	0.00
Spiked Amount 20.000	Range 70 - 130		Recovery = 92.90%			
36) d4-1,2-Dichloroethane	7.87	65	187003	18.58	ug/L	0.00
Spiked Amount 20.000	Range 70 - 130		Recovery = 92.90%			
54) d8-Toluene	10.74	98	451057	18.82	ug/L	0.00
Spiked Amount 20.000	Range 70 - 130		Recovery = 94.10%			
73) Bromofluorobenzene	14.89	95	174672	18.14	ug/L	0.00
Spiked Amount 20.000	Range 70 - 130		Recovery = 90.70%			
Target Compounds						Qvalue
2) Dichlorodifluoromethane	1.69	85	147051	17.21	ug/L	99
3) Chloromethane	1.93	50	155193	18.11	ug/L	97
4) Vinyl Chloride	2.01	62	156647	17.87	ug/L	97
5) Bromomethane	2.45	94	100699	22.19	ug/L	94
6) Chloroethane	2.54	64	78615	18.30	ug/L	99
7) t-Butyl alcohol (TBA)	3.87	59	68227	103.77	ug/L	96
8) Trichlorofluoromethane	2.85	101	292390	18.47	ug/L	100
9) Diethyl ether	3.25	74	49564	19.88	ug/L	94
10) 1,1,2-Trichlorotrifluoroet	3.42	101	130401	21.10	ug/L	99
11) Acrolein	3.40	56	94961	173.59	ug/L	98
12) Acetone	3.51	43	215786	98.18	ug/L	97
13) 1,1-Dichloroethene	3.65	96	97160	21.57	ug/L	97
14) Methyl iodide	4.09	142	211494	30.23	ug/L	98
15) Di-isopropyl ether (DIPE)	5.49	45	440299	19.82	ug/L	98
16) Methylene Chloride	4.41	84	102135	20.80	ug/L	96
17) Carbon Disulfide	4.36	76	334278	23.16	ug/L	100
18) Acrylonitrile	4.63	53	45711	21.96	ug/L	93
19) Methyl-tert-butyl ether (M	4.71	73	356626	20.96	ug/L	99
20) trans-1,2-Dichloroethene	4.88	61	220334	19.90	ug/L	99
21) 1,1-Dichloroethane	5.57	63	261630	20.71	ug/L	98
22) Vinyl acetate	5.63	43	280534	26.61	ug/L #	95
23) Methyl ethyl ketone	6.31	72	58731	109.29	ug/L	98
24) Ethyl t-butyl ether (ETBE)	6.13	59	458459	22.03	ug/L	97
25) 2,2-Dichloropropane	6.45	77	195633	20.15	ug/L	96
26) cis-1,2-Dichloroethene	6.53	96	156399	21.79	ug/L	96
27) t-Amyl methyl ether (TAME)	7.87	73	321088	22.05	ug/L	97

(#) = qualifier out of range (m) = manual integration

Quantitation Report (QT Reviewed)

Data File : D:\HPCHEM\1\DATA\052206C\C59026SI.D Vial: 6
Acq On : 22 May 2006 1:07 pm Operator:
Sample : 20 PPB Inst : Instr_C
Misc : 5000 Multiplr: 1.00
Integration Params: rteint.p
Quant Time: May 22 13:18 2006 Quant Results File: V805226C.RES

Quant Method : D:\HPCHEM\1\METHODS\V805226C.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Wed Apr 19 15:41:22 2006
Response via : Initial Calibration
DataAcq Meth : V805226C

	Compound	R.T.	QIon	Response	Conc Unit	Qvalue
29)	Chloroform	6.78	83	251002	21.06 ug/L	97
30)	Bromochloromethane	7.01	128	83164	21.85 ug/L	94
31)	Tetrahydrofuran	7.12	71	13852	23.85 ug/L	95
32)	1,1,1-Trichloroethane	7.37	97	227662	21.00 ug/L	97
34)	1,1-Dichloropropene	7.62	75	185921	21.55 ug/L	97
35)	Carbon Tetrachloride	7.75	119	198942	22.24 ug/L	95
37)	1,2-Dichloroethane	8.01	62	240228	20.42 ug/L	99
38)	Benzene	8.01	78	441081	21.46 ug/L	98
40)	Trichloroethene	8.94	95	154398	20.50 ug/L	96
41)	1,2-Dichloropropane	9.21	63	143246	20.21 ug/L	99
42)	Methylmethacrylate	9.37	69	91942	20.95 ug/L	97
43)	Methyl acetate	4.22	43	113579	17.55 ug/L	99
44)	Cyclohexane	7.38	56	235271	20.52 ug/L	96
45)	Bromodichloromethane	9.55	83	183028	19.76 ug/L	98
46)	Dibromomethane	9.62	93	115476	21.33 ug/L	99
47)	1,4-Dioxane	9.63	88	35110	524.17 ug/L	99
48)	Methylcyclohexane	9.03	83	197167	21.56 ug/L	95
50)	2-Chloroethylvinyl ether	10.07	63	48974	23.34 ug/L	99
51)	2-Hexanone	11.44	58	253739	107.48 ug/L	97
52)	Methyl isobutyl ketone	10.13	58	265744	109.16 ug/L	100
53)	cis-1,3-Dichloropropene	10.39	75	213977	20.30 ug/L	96
55)	Toluene	10.86	92	300584	21.16 ug/L	100
56)	trans-1,3-Dichloropropene	11.15	75	197060	19.64 ug/L	97
58)	1,1,2-Trichloroethane	11.38	83	98897	21.18 ug/L	93
59)	1,3-Dichloropropane	11.75	76	189115	20.71 ug/L	99
60)	Tetrachloroethene	11.83	166	178624	19.56 ug/L	97
61)	Dibromochloromethane	12.12	129	159300	20.24 ug/L	99
62)	1,2-Dibromoethane	12.42	107	167915	21.23 ug/L	95
64)	Chlorobenzene	13.07	112	395676	20.71 ug/L	97
65)	1,1,1,2-Tetrachloroethane	13.15	131	140023	20.67 ug/L	99
66)	Ethylbenzene	13.17	91	569563	20.63 ug/L	99
67)	m,p-Xylene	13.29	106	453702	40.99 ug/L	97
68)	o-Xylene	13.95	106	217154	20.27 ug/L	97
69)	Styrene	14.01	104	386629	20.52 ug/L	94
70)	Bromoform	14.49	173	120862	20.56 ug/L	96
71)	Isopropylbenzene	14.52	105	583100	20.51 ug/L	99
72)	1,1,2,2-Tetrachloroethane	14.78	83	152248	20.48 ug/L	97
74)	1,2,3-Trichloropropane	15.01	75	110342	19.98 ug/L	97
75)	trans-1,4-Dichloro-2-buten	15.14	89	45099	19.61 ug/L	98
76)	n-Propylbenzene	15.15	91	647161	20.33 ug/L	100
77)	Bromobenzene	15.17	156	163319	19.95 ug/L	97

(#) = qualifier out of range (m) = manual integration

Quantitation Report (QT Reviewed)

Data File : D:\HPCHEM\1\DATA\052206C\C59026SI.D Vial: 6
Acq On : 22 May 2006 1:07 pm Operator:
Sample : 20 PPB Inst : Instr_C
Misc : 5000 Multiplr: 1.00
Integration Params: rteint.p
Quant Time: May 22 13:18 2006 Quant Results File: V805226C.RES

Quant Method : D:\HPCHEM\1\METHODS\V805226C.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Wed Apr 19 15:41:22 2006
Response via : Initial Calibration
DataAcq Meth : V805226C

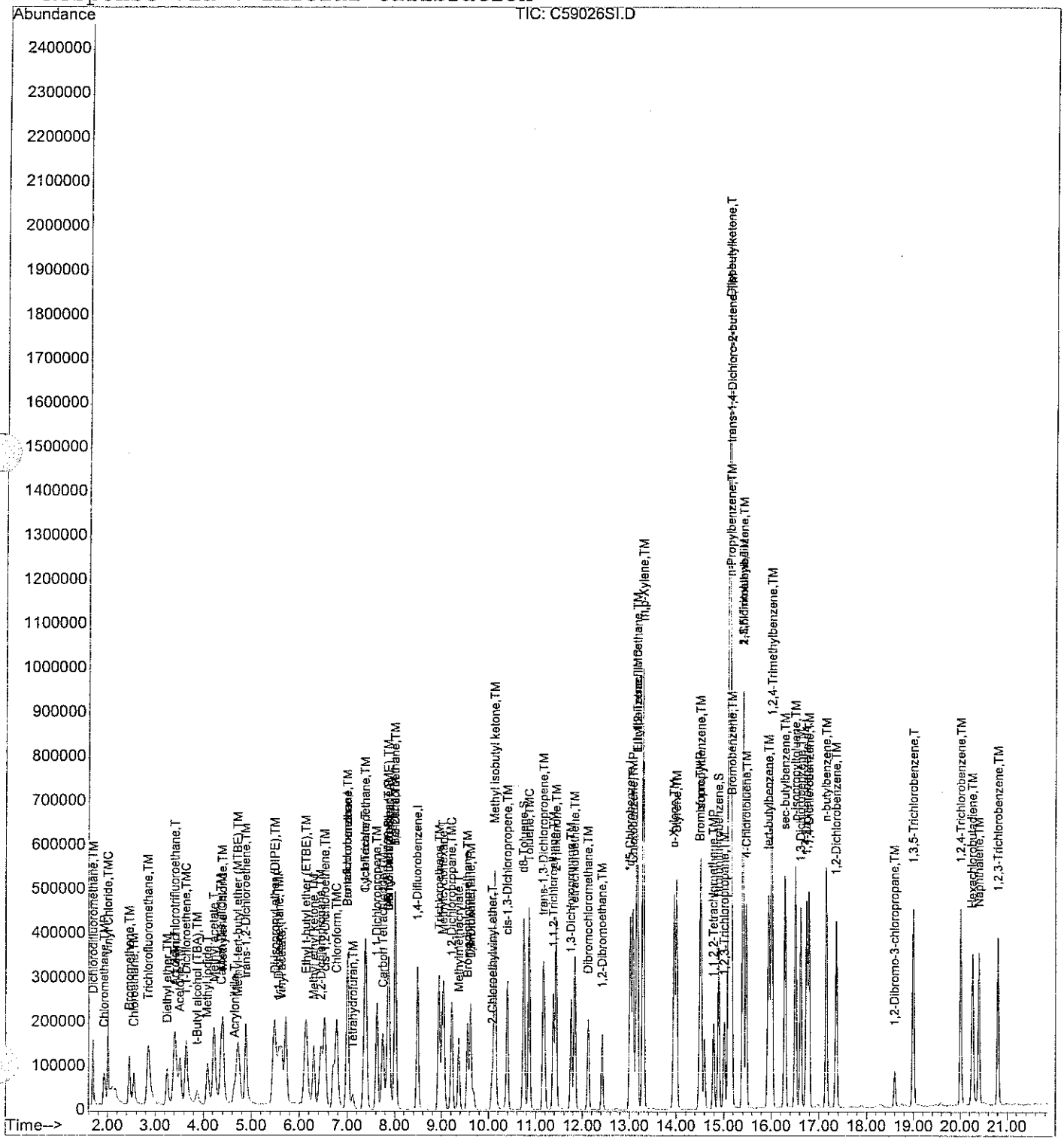
	Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
78)	1,3,5-Trimethylbenzene	15.40	105	465575	19.84	ug/L	99
79)	2-Chlorotoluene	15.39	91	433946	19.78	ug/L	99
80)	4-Chlorotoluene	15.48	91	385985	19.16	ug/L	99
81)	tert-butylbenzene	15.94	119	404035	19.48	ug/L	98
82)	1,2,4-Trimethylbenzene	16.00	105	455297	19.08	ug/L	99
83)	Diisobutylketone	15.11	57	1675468	100.26	ug/L	96
84)	sec-butylbenzene	16.28	105	627665	19.65	ug/L	97
85)	p-isopropyltoluene	16.51	119	502149	19.58	ug/L	96
86)	1,3-Dichlorobenzene	16.61	146	270612	19.56	ug/L	97
88)	1,4-Dichlorobenzene	16.78	146	280208	20.76	ug/L	98
89)	n-butylbenzene	17.15	91	467207	20.66	ug/L	99
90)	1,2-Dichlorobenzene	17.35	146	261172	20.97	ug/L	98
92)	1,2-Dibromo-3-chloropropan	18.60	75	25438	19.12	ug/L	83
94)	1,2,4-Trichlorobenzene	20.01	180	212907	19.90	ug/L	97
95)	Hexachlorobutadiene	20.25	225	104878	19.19	ug/L	98
96)	Naphthalene	20.39	128	388778	19.92	ug/L	100
97)	1,2,3-Trichlorobenzene	20.81	180	187174	19.29	ug/L	98
98)	1,3,5-Trichlorobenzene	19.00	180	235701	19.61	ug/L	100

(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : D:\HPCHEM\1\DATA\052206C\C59026SI.D Vial: 6
Acq On : 22 May 2006 1:07 pm Operator:
Sample : 20 PPB Inst : Instr_C
Misc : 5000 Multiplr: 1.00
MS Integration Params: rteint.p
Quant Time: May 22 13:18 2006 Quant Results File: V805226C.RES

Method : D:\HPCHEM\1\METHODS\V805226C.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Mon May 22 14:18:09 2006
Response via : Initial Calibration



Quantitation Report (Not Reviewed)

Data File : D:\HPCHEM\1\DATA\052206C\C59027SI.D Vial: 7
Acq On : 22 May 2006 1:07 pm Operator:
Sample : 50 PPB Inst : Instr_C
Misc : 5000 Multiplr: 1.00
Integration Params: rteint.p
Quant Time: May 22 13:29 2006 Quant Results File: V805226C.RES

Quant Method : D:\HPCHEM\1\METHODS\V805226C.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Wed Apr 19 15:41:22 2006
Response via : Initial Calibration
DataAcq Meth : V805226C

MT 5/23/06

5.27.06

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.00	168	239425	20.00	ug/L	-0.02
39) 1,4-Difluorobenzene	8.49	114	405568	20.00	ug/L	0.00
63) *d5-Chlorobenzene	13.01	117	362863	20.00	ug/L	0.00
87) 1,4-Dichlorobenzene-d4	16.74	152	178724	20.00	ug/L	0.00
System Monitoring Compounds						
28) DB Holder #28	7.86	65	436774	45.90	ug/L	0.00
Spiked Amount 20.000	Range 70 - 130		Recovery = 229.50%#			
36) d4-1,2-Dichloroethane	7.86	65	436774	45.90	ug/L	0.00
Spiked Amount 20.000	Range 70 - 130		Recovery = 229.50%#			
54) d8-Toluene	10.74	98	1060126	46.71	ug/L	0.00
Spiked Amount 20.000	Range 70 - 130		Recovery = 233.55%#			
73) Bromofluorobenzene	14.90	95	424602	46.05	ug/L	0.00
Spiked Amount 20.000	Range 70 - 130		Recovery = 230.25%#			
Target Compounds						Qvalue
2) Dichlorodifluoromethane	1.68	85	354601	43.91	ug/L	99
3) Chloromethane	1.95	50	405998	50.12	ug/L	99
4) Vinyl Chloride	2.01	62	350028	42.23	ug/L	99
5) Bromomethane	2.43	94	254219	59.27	ug/L	100
6) Chloroethane	2.52	64	187263	46.10	ug/L	100
7) t-Butyl alcohol (TBA)	3.93	59	155913	250.84	ug/L	99
8) Trichlorofluoromethane	2.83	101	701434	46.86	ug/L	97
9) Diethyl ether	3.26	74	113844	48.30	ug/L	98
10) 1,1,2-Trichlorotrifluoroet	3.42	101	305908	52.35	ug/L	98
11) Acrolein	3.41	56	231651	447.93	ug/L	99
12) Acetone	3.54	43	527816	254.02	ug/L	99
13) 1,1-Dichloroethene	3.63	96	225245	52.91	ug/L	97
14) Methyl iodide	4.08	142	426910	64.54	ug/L	97
15) Di-isopropyl ether (DIPE)	5.49	45	1033311	49.19	ug/L	98
16) Methylene Chloride	4.39	84	234215	50.45	ug/L	98
17) Carbon Disulfide	4.34	76	803024	58.86	ug/L	100
18) Acrylonitrile	4.64	53	106680	54.22	ug/L	95
19) Methyl-tert-butyl ether (M	4.72	73	839984	52.23	ug/L	99
20) trans-1,2-Dichloroethene	4.88	61	538766	51.47	ug/L	99
21) 1,1-Dichloroethane	5.57	63	628076	52.60	ug/L	98
22) Vinyl acetate	5.63	43	706802	70.93	ug/L #	95
23) Methyl ethyl ketone	6.31	72	135957	267.62	ug/L	97
24) Ethyl t-butyl ether (ETBE)	6.14	59	1097739	55.79	ug/L	97
25) 2,2-Dichloropropane	6.45	77	442213	48.18	ug/L	97
26) cis-1,2-Dichloroethene	6.52	96	359138	52.92	ug/L	99
27) t-Amyl methyl ether (TAME)	7.86	73	749762	54.48	ug/L	98

(#) = qualifier out of range (m) = manual integration

Quantitation Report (Not Reviewed)

Data File : D:\HPCHEM\1\DATA\052206C\C59027SI.D Vial: 7
Acq On : 22 May 2006 1:07 pm Operator:
Sample : 50 PPB Inst : Instr_C
Misc : 5000 Multiplr: 1.00
Integration Params: rteint.p
Quant Time: May 22 13:29 2006 Quant Results File: V805226C.RES

Quant Method : D:\HPCHEM\1\METHODS\V805226C.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Wed Apr 19 15:41:22 2006
Response via : Initial Calibration
DataAcq Meth : V805226C

	Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
29)	Chloroform	6.78	83	591345	52.49	ug/L	96
30)	Bromochloromethane	7.01	128	189192	52.57	ug/L	94
31)	Tetrahydrofuran	7.12	71	31754	57.83	ug/L	95
32)	1,1,1-Trichloroethane	7.37	97	541473	52.82	ug/L	99
34)	1,1-Dichloropropene	7.62	75	433749	53.18	ug/L	98
35)	Carbon Tetrachloride	7.75	119	479255	56.68	ug/L	99
37)	1,2-Dichloroethane	8.00	62	566654	50.96	ug/L	99
38)	Benzene	8.00	78	988625	50.88	ug/L	99
40)	Trichloroethene	8.95	95	366650	51.41	ug/L	99
41)	1,2-Dichloropropane	9.22	63	334498	49.82	ug/L	100
42)	Methylmethacrylate	9.36	69	203286	48.91	ug/L	99
43)	Methyl acetate	4.22	43	247012	40.30	ug/L	98
44)	Cyclohexane	7.38	56	542037	49.92	ug/L	97
45)	Bromodichloromethane	9.56	83	447280	50.99	ug/L	97
46)	Dibromomethane	9.62	93	262118	51.12	ug/L	97
47)	1,4-Dioxane	9.64	88	89518	1411.16	ug/L	99
48)	Methylcyclohexane	9.03	83	461180	53.26	ug/L	96
50)	2-Chloroethylvinyl ether	10.07	63	105001	52.85	ug/L	98
51)	2-Hexanone	11.43	58	606572	271.28	ug/L	96
52)	Methyl isobutyl ketone	10.13	58	598298	259.50	ug/L	99
53)	cis-1,3-Dichloropropene	10.39	75	512948	51.37	ug/L	97
55)	Toluene	10.85	92	694908	51.65	ug/L	99
56)	trans-1,3-Dichloropropene	11.15	75	470779	49.54	ug/L	100
58)	1,1,2-Trichloroethane	11.38	83	231972	52.45	ug/L	94
59)	1,3-Dichloropropane	11.76	76	441015	51.00	ug/L	99
60)	Tetrachloroethene	11.84	166	418087	48.33	ug/L	100
61)	Dibromochloromethane	12.12	129	397370	53.30	ug/L	100
62)	1,2-Dibromoethane	12.42	107	388893	51.93	ug/L	100
64)	Chlorobenzene	13.07	112	937166	51.23	ug/L	99
65)	1,1,1,2-Tetrachloroethane	13.15	131	341997	52.72	ug/L	99
66)	Ethylbenzene	13.17	91	1364256	51.60	ug/L	100
67)	m,p-Xylene	13.29	106	1048205	98.90	ug/L	99
68)	o-Xylene	13.96	106	527268	51.40	ug/L	98
69)	Styrene	14.01	104	916941	50.84	ug/L	96
70)	Bromoform	14.49	173	323046	57.39	ug/L	96
71)	Isopropylbenzene	14.52	105	1397219	51.33	ug/L	100
72)	1,1,2,2-Tetrachloroethane	14.79	83	367401	51.62	ug/L	99
74)	1,2,3-Trichloropropane	15.00	75	270397	51.14	ug/L	98
75)	trans-1,4-Dichloro-2-buten	15.14	89	109763	49.85	ug/L	99
76)	n-Propylbenzene	15.14	91	1505547	49.40	ug/L	97
77)	Bromobenzene	15.17	156	403583	51.48	ug/L	96

(#) = qualifier out of range (m) = manual integration

Quantitation Report (Not Reviewed)

Data File : D:\HPCHEM\1\DATA\052206C\C59027SI.D Vial: 7
Acq On : 22 May 2006 1:07 pm Operator:
Sample : 50 PPB Inst : Instr_C
Misc : 5000 Multiplr: 1.00
Integration Params: rteint.p
Quant Time: May 22 13:29 2006 Quant Results File: V805226C.RES

Quant Method : D:\HPCHEM\1\METHODS\V805226C.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Wed Apr 19 15:41:22 2006
Response via : Initial Calibration
DataAcq Meth : V805226C

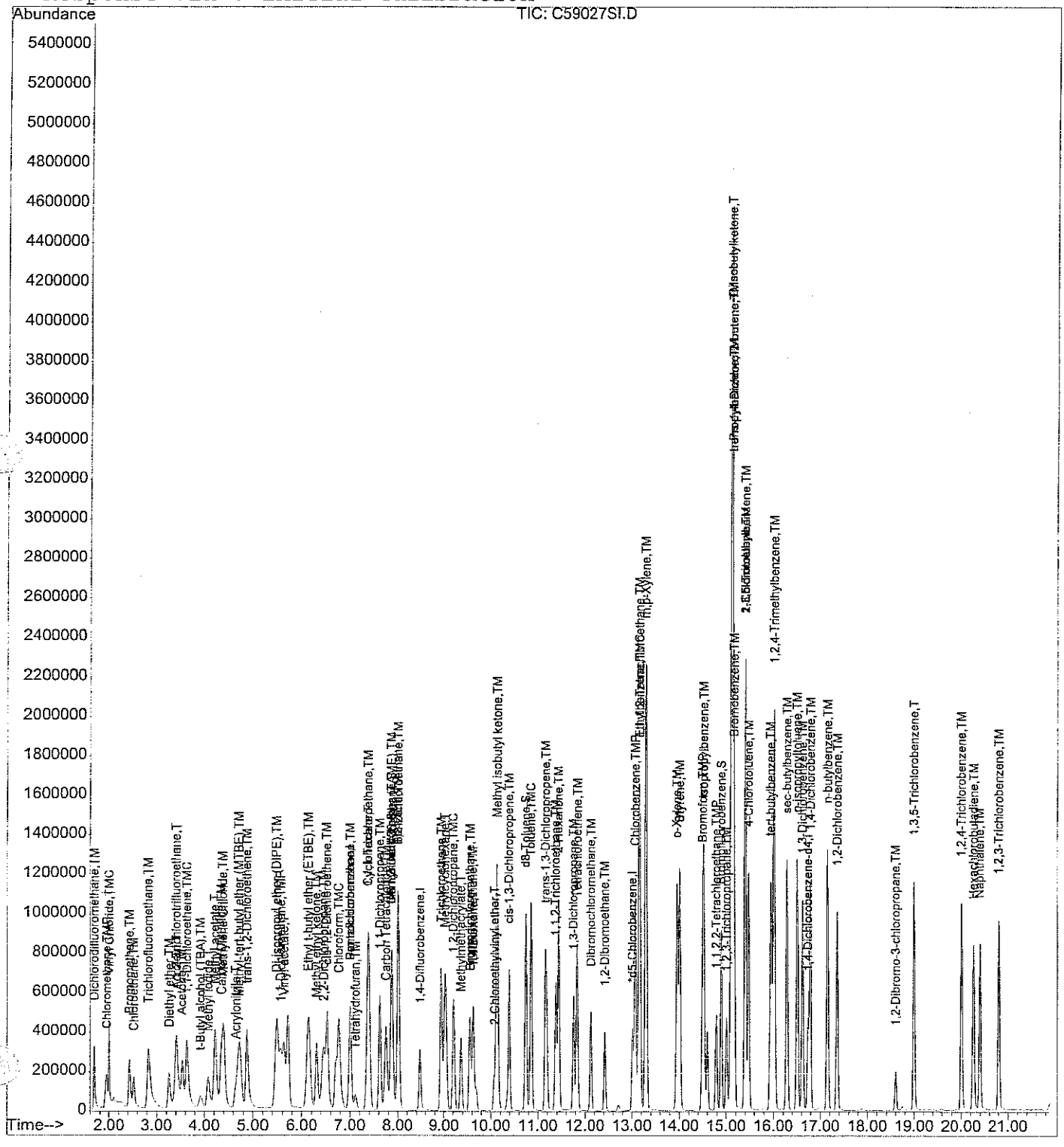
	Compound	R.T.	QIon	Response	Conc	Unit		Qvalue
78)	1,3,5-Trimethylbenzene	15.40	105	1128527	50.24	ug/L		100
79)	2-Chlorotoluene	15.40	91	1030115	49.04	ug/L		99
80)	4-Chlorotoluene	15.48	91	974529	50.52	ug/L		99
81)	tert-butylbenzene	15.95	119	1006837	50.71	ug/L		97
82)	1,2,4-Trimethylbenzene	16.01	105	1155132	50.56	ug/L		99
83)	Diisobutylketone	15.11	57	3901432	243.84	ug/L		98
84)	sec-butylbenzene	16.29	105	1544039	50.48	ug/L		97
85)	p-isopropyltoluene	16.50	119	1223522	49.82	ug/L		98
86)	1,3-Dichlorobenzene	16.62	146	662344	50.01	ug/L		97
88)	1,4-Dichlorobenzene	16.79	146	695137	52.05	ug/L		99
89)	n-butylbenzene	17.15	91	1138812	50.89	ug/L		98
90)	1,2-Dichlorobenzene	17.36	146	626664	50.84	ug/L		100
92)	1,2-Dibromo-3-chloropropan	18.60	75	67039	50.93	ug/L	#	75
94)	1,2,4-Trichlorobenzene	20.01	180	533358	50.39	ug/L		96
95)	Hexachlorobutadiene	20.26	225	272102	50.30	ug/L		97
96)	Naphthalene	20.40	128	1021768	52.91	ug/L		100
97)	1,2,3-Trichlorobenzene	20.80	180	492108	51.24	ug/L		100
98)	1,3,5-Trichlorobenzene	19.00	180	599211	50.38	ug/L		100

(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : D:\HPCHEM\1\DATA\052206C\C59027SI.D Vial: 7
Acq On : 22 May 2006 1:07 pm Operator:
Sample : 50 PPB Inst : Instr_C
Misc : 5000 Multiplr: 1.00
MS Integration Params: rteint.p
Quant Time: May 22 13:29 2006 Quant Results File: V805226C.RES

Method : D:\HPCHEM\1\METHODS\V805226C.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Mon May 22 14:18:09 2006
Response via : Initial Calibration



Quantitation Report (QT Reviewed)

Data File : D:\HPCHEM\1\DATA\052206C\C59028SI.D Vial: 8
Acq On : 22 May 2006 1:50 pm Operator:
Sample : 100 PPB Inst : Instr_C
Misc : 5000 Multiplr: 1.00
Integration Params: rteint.p
Quant Time: May 22 14:03 2006 Quant Results File: V805226C.RES

Quant Method : D:\HPCHEM\1\METHODS\V805226C.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Wed Apr 19 15:41:22 2006
Response via : Initial Calibration
DataAcq Meth : V805226C

MT 5/23/06 5.23.06

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	6.98	168	221068	20.00	ug/L	-0.03
39) 1,4-Difluorobenzene	8.47	114	392598	20.00	ug/L	-0.02
63) *d5-Chlorobenzene	13.02	117	366401 ✓	20.00	ug/L	0.00
87) 1,4-Dichlorobenzene-d4	16.74	152	173234	20.00	ug/L	0.00
System Monitoring Compounds						
28) DB Holder #28	7.86	65	868591	98.86	ug/L	-0.01
Spiked Amount 20.000	Range 70 - 130		Recovery = 494.30%#			
36) d4-1,2-Dichloroethane	7.86	65	868591	98.86	ug/L	-0.01
Spiked Amount 20.000	Range 70 - 130		Recovery = 494.30%#			
54) d8-Toluene	10.74	98	2060815	93.80	ug/L	0.00
Spiked Amount 20.000	Range 70 - 130		Recovery = 469.00%#			
73) Bromofluorobenzene	14.89	95	828520	89.00	ug/L	0.00
Spiked Amount 20.000	Range 70 - 130		Recovery = 445.00%#			
Target Compounds						Qvalue
2) Dichlorodifluoromethane	1.68	85	683055	91.61	ug/L	99
3) Chloromethane	1.93	50	850952	113.77	ug/L	96
4) Vinyl Chloride	1.99	62	687475	89.84	ug/L	100
5) Bromomethane	2.43	94	127880	32.29	ug/L	98
6) Chloroethane	2.49	64	398203	106.17	ug/L	99
7) t-Butyl alcohol (TBA)	4.02	59	314571m	548.13	ug/L	
8) Trichlorofluoromethane	2.79	101	1387906	100.42	ug/L	100
9) Diethyl ether	3.26	74	221250	101.67	ug/L	99
10) 1,1,2-Trichlorotrifluoroet	3.37	101	594837	110.25	ug/L	99
11) Acrolein	3.41	56	279728	585.80	ug/L	99
12) Acetone	3.56	43	1013377	528.20	ug/L	100
13) 1,1-Dichloroethene	3.61	96	437984	111.42	ug/L	96
14) Methyl iodide	4.05	142	898075	147.04	ug/L	97
15) Di-isopropyl ether (DIPE)	5.49	45	2025383	104.43	ug/L	97
16) Methylene Chloride	4.38	84	453094	105.69	ug/L	97
17) Carbon Disulfide	4.32	76	1573923	124.94	ug/L	100
18) Acrylonitrile	4.65	53	215735	118.75	ug/L	95
19) Methyl-tert-butyl ether (M	4.73	73	1661291	111.87	ug/L	99
20) trans-1,2-Dichloroethene	4.86	61	1047766	108.40	ug/L	99
21) 1,1-Dichloroethane	5.56	63	1253634	113.70	ug/L	99
22) Vinyl acetate	5.63	43	1285389	139.70	ug/L #	95
23) Methyl ethyl ketone	6.32	72	275380	587.07	ug/L	100
24) Ethyl t-butyl ether (ETBE)	6.15	59	2201212	121.17	ug/L	97
25) 2,2-Dichloropropane	6.44	77	1344317	158.64	ug/L	88
26) cis-1,2-Dichloroethene	6.51	96	644651	102.88	ug/L #	79
27) t-Amyl methyl ether (TAME)	7.87	73	1437720	113.14	ug/L	98

(#) = qualifier out of range (m) = manual integration

Quantitation Report (QT Reviewed)

Data File : D:\HPCHEM\1\DATA\052206C\C59028SI.D Vial: 8
Acq On : 22 May 2006 1:50 pm Operator:
Sample : 100 PPB Inst : Instr_C
Misc : 5000 Multiplr: 1.00
Integration Params: rteint.p
Quant Time: May 22 14:03 2006 Quant Results File: V805226C.RES

Quant Method : D:\HPCHEM\1\METHODS\V805226C.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Wed Apr 19 15:41:22 2006
Response via : Initial Calibration
DataAcq Meth : V805226C

	Compound	R.T.	QIon	Response	Conc Unit		Qvalue
29)	Chloroform	6.77	83	1562921	150.25 ug/L		94
30)	Bromochloromethane	6.99	128	360876	108.60 ug/L	#	85
31)	Tetrahydrofuran	7.13	71	62456	123.18 ug/L		95
32)	1,1,1-Trichloroethane	7.36	97	1082377	114.36 ug/L		98
34)	1,1-Dichloropropene	7.60	75	834734	110.83 ug/L		99
35)	Carbon Tetrachloride	7.73	119	979753	125.49 ug/L		99
37)	1,2-Dichloroethane	8.00	62	1155448	112.53 ug/L		99
38)	Benzene	8.00	78	1897298	105.76 ug/L		100
40)	Trichloroethene	8.94	95	702470	101.75 ug/L		98
41)	1,2-Dichloropropane	9.21	63	636013	97.86 ug/L		100
42)	Methylmethacrylate	9.37	69	412112	102.43 ug/L		99
43)	Methyl acetate	4.23	43	501763	84.57 ug/L		99
44)	Cyclohexane	7.36	56	984448	93.66 ug/L		98
45)	Bromodichloromethane	9.56	83	911003	107.28 ug/L		98
46)	Dibromomethane	9.62	93	523851	105.54 ug/L		98
47)	1,4-Dioxane	9.66	88	168813	2749.07 ug/L		99
48)	Methylcyclohexane	9.02	83	888275	105.97 ug/L		96
50)	2-Chloroethylvinyl ether	10.07	63	334036	173.67 ug/L		100
51)	2-Hexanone	11.45	58	1144173	528.63 ug/L		97
52)	Methyl isobutyl ketone	10.14	58	1153904	517.02 ug/L		99
53)	cis-1,3-Dichloropropene	10.40	75	1005518	104.03 ug/L		97
55)	Toluene	10.86	92	1353323	103.91 ug/L		98
56)	trans-1,3-Dichloropropene	11.15	75	967583	105.19 ug/L		99
58)	1,1,2-Trichloroethane	11.38	83	447275	104.48 ug/L		96
59)	1,3-Dichloropropane	11.76	76	866204	103.49 ug/L		98
60)	Tetrachloroethene	11.84	166	840060	100.32 ug/L		98
61)	Dibromochloromethane	12.12	129	803386	111.32 ug/L		98
62)	1,2-Dibromoethane	12.43	107	792463	109.31 ug/L		99
64)	Chlorobenzene	13.07	112	1830001	99.08 ug/L		98
65)	1,1,1,2-Tetrachloroethane	13.15	131	681347	104.02 ug/L		97
66)	Ethylbenzene	13.18	91	2569943	96.27 ug/L		99
67)	m,p-Xylene	13.29	106	1930987	180.44 ug/L		91
68)	o-Xylene	13.96	106	1005184	97.04 ug/L		95
69)	Styrene	14.01	104	1795152	98.57 ug/L		98
70)	Bromoform	14.50	173	644889	113.45 ug/L		96
71)	Isopropylbenzene	14.52	105	2668677	97.10 ug/L		99
72)	1,1,2,2-Tetrachloroethane	14.78	83	707778	98.49 ug/L		98
74)	1,2,3-Trichloropropane	15.01	75	516718	96.79 ug/L		97
75)	trans-1,4-Dichloro-2-buten	15.14	89	209418	94.20 ug/L		97
76)	n-Propylbenzene	15.15	91	2671037	86.79 ug/L		97
77)	Bromobenzene	15.18	156	760045	96.01 ug/L		97

(#) = qualifier out of range (m) = manual integration

Quantitation Report (QT Reviewed)

Data File : D:\HPCHEM\1\DATA\052206C\C59028SI.D Vial: 8
Acq On : 22 May 2006 1:50 pm Operator:
Sample : 100 PPB Inst : Instr_C
Misc : 5000 Multiplr: 1.00
Integration Params: rteint.p
Quant Time: May 22 14:03 2006 Quant Results File: V805226C.RES

Quant Method : D:\HPCHEM\1\METHODS\V805226C.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Wed Apr 19 15:41:22 2006
Response via : Initial Calibration
DataAcq Meth : V805226C

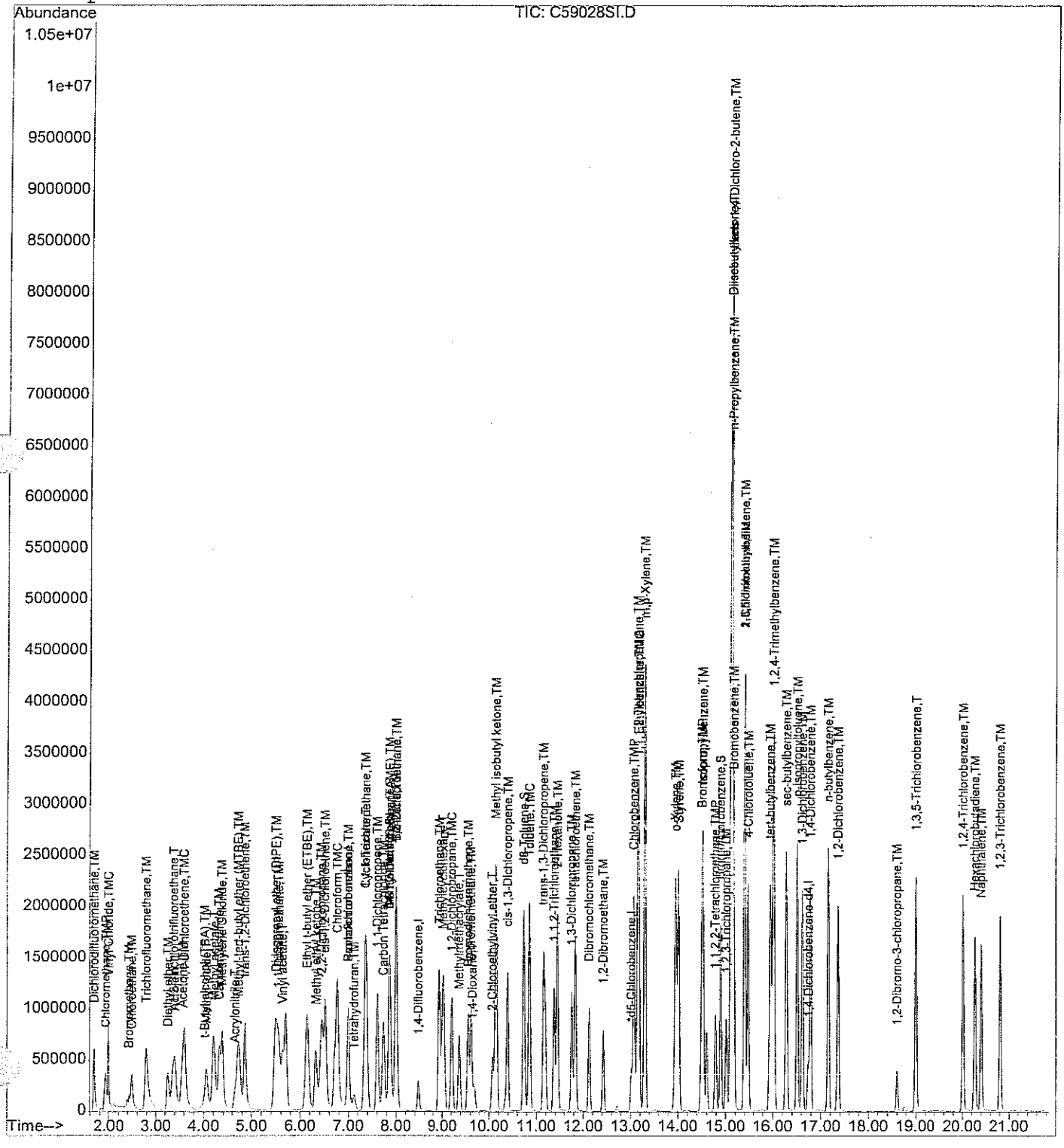
	Compound	R.T.	QIon	Response	Conc Unit	Qvalue
78)	1,3,5-Trimethylbenzene	15.40	105	2136796	94.21 ug/L	98
79)	2-Chlorotoluene	15.39	91	1986483	93.66 ug/L	97
80)	4-Chlorotoluene	15.48	91	1848978	94.93 ug/L	97
81)	tert-butylbenzene	15.95	119	1868398	93.19 ug/L	97
82)	1,2,4-Trimethylbenzene	16.01	105	2179132	94.46 ug/L	97
83)	Diisobutylketone	15.12	57	7102212	439.61 ug/L	96
84)	sec-butylbenzene	16.28	105	3013729	97.59 ug/L	97
85)	p-isopropyltoluene	16.51	119	2361350	95.22 ug/L	99
86)	1,3-Dichlorobenzene	16.62	146	1288202	96.32 ug/L	98
88)	1,4-Dichlorobenzene	16.79	146	1297093	100.21 ug/L	99
89)	n-butylbenzene	17.15	91	2177244	100.38 ug/L	99
90)	1,2-Dichlorobenzene	17.36	146	1192521	99.81 ug/L	99
92)	1,2-Dibromo-3-chloropropan	18.60	75	129412	101.42 ug/L	84
94)	1,2,4-Trichlorobenzene	20.01	180	1053074	102.64 ug/L	97
95)	Hexachlorobutadiene	20.26	225	538018	102.61 ug/L	98
96)	Naphthalene	20.39	128	1934665	103.36 ug/L	100
97)	1,2,3-Trichlorobenzene	20.81	180	950271	102.09 ug/L	99
98)	1,3,5-Trichlorobenzene	19.00	180	1188399	103.07 ug/L	98

(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : D:\HPCHEM\1\DATA\052206C\C59028SI.D Vial: 8
Acq On : 22 May 2006 1:50 pm Operator:
Sample : 100 PPB Inst : Instr_C
Misc : 5000 Multiplr: 1.00
MS Integration Params: rteint.p
Quant Time: May 22 14:03 2006 Quant Results File: V805226C.RES

Method : D:\HPCHEM\1\METHODS\V805226C.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Mon May 22 14:18:09 2006
Response via : Initial Calibration

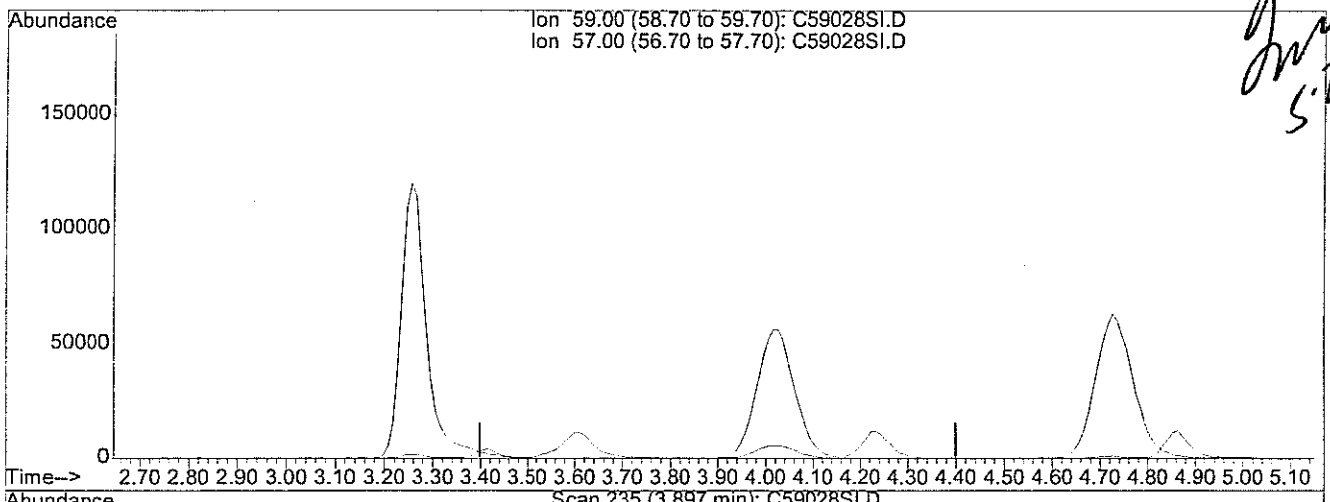


Quantitation Report (Qedit)

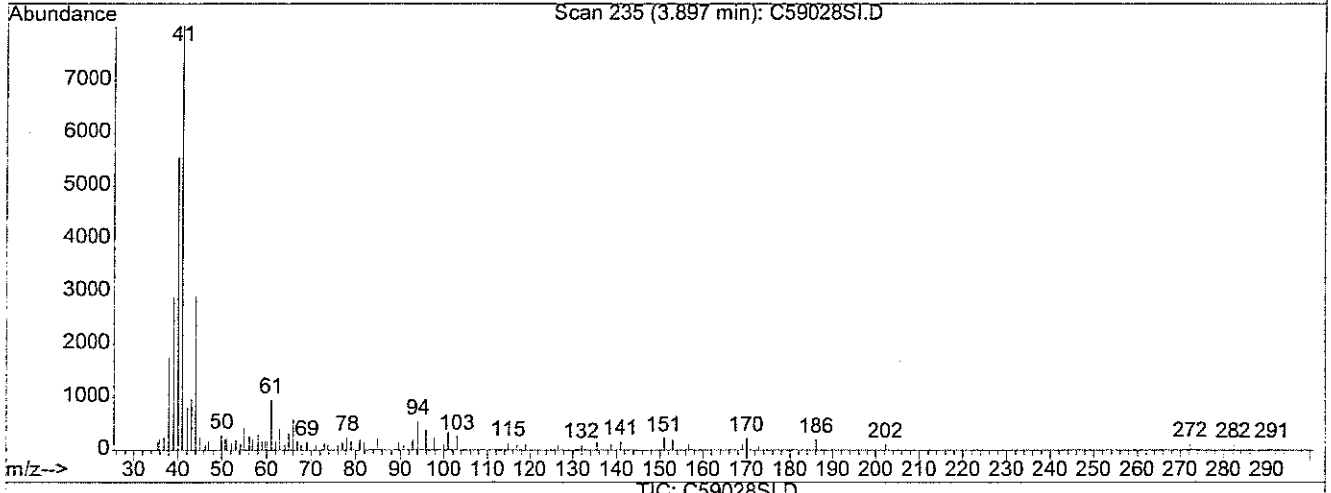
Data File : D:\HPCHEM\1\DATA\052206C\C59028SI.D Vial: 8
Acq On : 22 May 2006 1:50 pm Operator:
Sample : 100 PPB Inst : Instr_C
Misc : 5000 Multiplr: 1.00
MS Integration Params: rteint.p
Quant Time: May 22 14:03 2006 Quant Results File: temp.res

Method : D:\HPCHEM\1\METHODS\V805226C.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Wed Apr 19 15:41:22 2006
Response via : Multiple Level Calibration

MT 5/23/06
missed peak



5.23.06



TIC: C59028SI.D

(7) t-Butyl alcohol (TBA) (TM)

3.90min 0.00ug/L

response 0

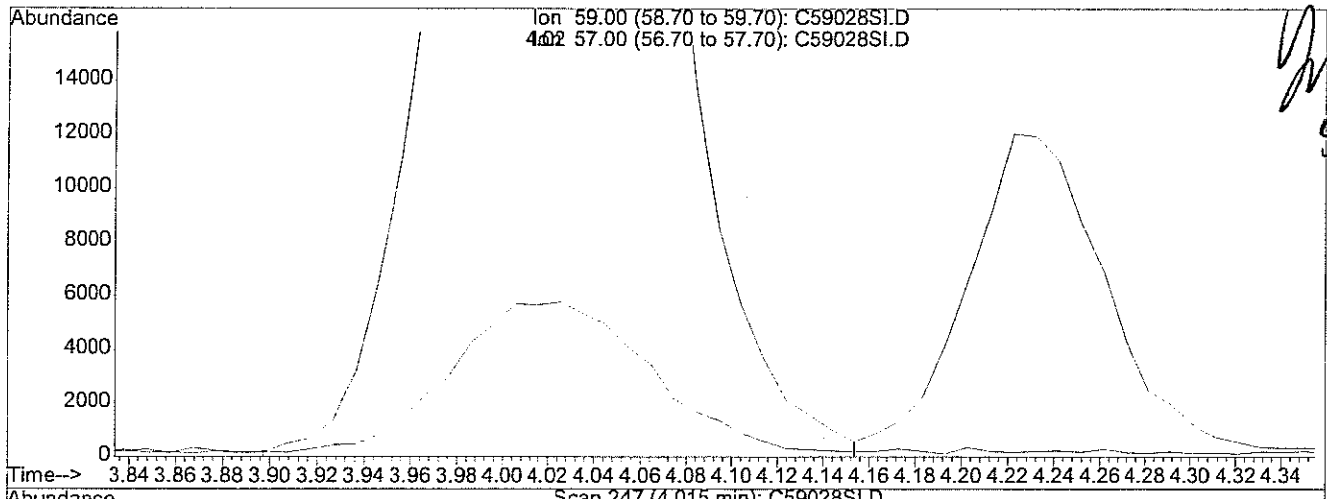
Ion	Exp%	Act%
59.00	100	0.00
57.00	11.60	0.00#
0.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (Qedit)

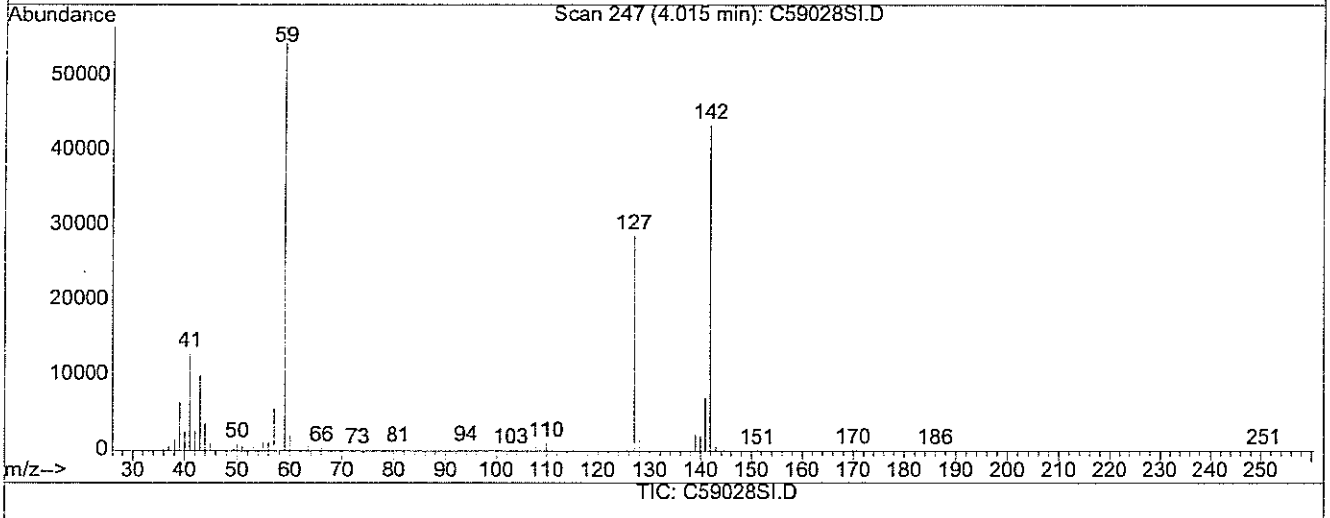
Data File : D:\HPCHEM\1\DATA\052206C\C59028SI.D Vial: 8
Acq On : 22 May 2006 1:50 pm Operator:
Sample : 100 PPB Inst : Instr_C
Misc : 5000 Multiplr: 1.00
MS Integration Params: rteint.p
Quant Time: May 22 14:03 2006 Quant Results File: temp.res

Method : D:\HPCHEM\1\METHODS\V805226C.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Wed Apr 19 15:41:22 2006
Response via : Multiple Level Calibration

MT 5/23/06



5.23.06



TIC: C59028SI.D

(7) t-Butyl alcohol (TBA) (TM)

4.02min 548.13ug/L m

response 314571

Ion	Exp%	Act%
59.00	100	100
57.00	11.60	0.00#
0.00	0.00	0.00
0.00	0.00	0.00

Acq. Method File	**V805266B**
Level 1 File Name	B36408SI.D
Level 2 File Name	B36409SI.D
Level 3 File Name	B36410SI.D
Level 4 File Name	B36411SI.D
Level 5 File Name	B36412SI.D
Level 6 File Name	B36413SI.D

6-30-06 SL ✓ 6-30-06

Name	Level 1 RF	Level 2 RF	Level 3 RF	Level 4 RF	Level 5 RF	Level 6 RF	AVG RF	RF Rel. STDEV
Dichlorodifluoromethane	0.695	0.687	0.694	0.688	0.676	0.656	0.683	2.17%
Chloromethane	1.382	1.290	1.129	1.164	1.097		1.212	9.88%
Vinyl Chloride	0.882	0.791	0.585	0.553	0.512		0.665	24.40% OK CCC
Bromomethane	0.343	0.262	0.320	0.447	0.583	0.591	0.424	32.86%
Chloroethane	0.763	0.631	0.618	0.603	0.578	0.540	0.622	12.24%
t-Butyl alcohol (TBA)	0.050	0.056	0.065	0.073	0.068		0.062	14.99%
Trichlorofluoromethane	1.190	1.175	1.231	1.180	1.168	1.125	1.178	2.90%
Diethyl ether	0.439	0.468	0.450	0.522	0.489	0.437	0.467	7.07%
1,1,2-Trichlorotrifluoroethane	0.504	0.499	0.518	0.516	0.516	0.510	0.511	1.52%
Acrolein	0.323	0.367	0.361	0.387	0.394	0.375	0.368	6.86%
Acetone	0.331	0.297	0.299	0.307	0.304	0.281	0.303	5.34%
1,1-Dichloroethene	0.471	0.449	0.461	0.480	0.488	0.493	0.474	3.54%
Methyl iodide	0.180	0.245	0.400	0.572	0.681	0.697	0.462	99.95% NT
Di-isopropyl ether (DIPE)	2.597	2.385	2.445	2.466	2.505	2.509	2.484	2.88%
Methylene Chloride	0.749	0.674	0.630	0.629	0.621		0.660	8.10%
Carbon Disulfide	1.904	1.779	1.798	1.869	1.939	1.960	1.875	3.95%
Acrylonitrile	0.283	0.319	0.305	0.337	0.337	0.325	0.318	6.53%
Methyl-tert-butyl ether (MTBE)	1.763	1.701	1.750	1.786	1.778	1.754	1.755	1.70%
trans-1,2-Dichloroethene	0.606	0.539	0.525	0.548	0.549	0.555	0.553	5.04%
1,1-Dichloroethane	1.145	1.095	1.081	1.090	1.114	1.124	1.108	2.17%
Vinyl acetate	0.030	0.060	0.079	0.104	0.117	0.116	0.084	99.96% NT
Methyl ethyl ketone	0.429	0.436	0.447	0.486	0.474	0.443	0.453	4.92%
Ethyl t-butyl ether (ETBE)	2.027	2.007	1.934	2.042	2.033	2.056	2.017	2.16%
2,2-Dichloropropane	0.696	0.631	0.644	0.719	0.765	0.767	0.703	8.28%
cis-1,2-Dichloroethene	0.593	0.566	0.566	0.586	0.598	0.599	0.585	2.60%
t-Amyl methyl ether (TAME)	1.688	1.637	1.655	1.727	1.733	1.724	1.694	2.40%
DB HOLDER #28	0.829	0.828	0.820	0.806	0.759	0.768	0.802	3.84%
Chloroform	1.188	1.031	1.055	1.065	1.066	1.067	1.079	5.13%
Bromochloromethane	0.259	0.259	0.263	0.263	0.263	0.261	0.261	0.77%
Tetrahydrofuran	0.092	0.084	0.092	0.092	0.086		0.089	4.22%
1,1,1-Trichloroethane	0.794	0.729	0.741	0.762	0.769	0.786	0.763	3.27%
1,1-Dichloropropene	0.879	0.813	0.799	0.817	0.806	0.809	0.821	3.54%
Carbon Tetrachloride	0.465	0.432	0.479	0.507	0.525	0.545	0.492	8.46%
d4-1,2-Dichloroethane	0.829	0.828	0.820	0.806	0.759	0.768	0.802	3.84%
1,2-Dichloroethane	1.129	1.043	1.005	0.979	0.966	0.949	1.012	6.54%
Benzene	2.566	2.235	2.216	2.282	2.272	2.322	2.316	5.53%
Trichloroethene	0.382	0.361	0.342	0.358	0.371	0.376	0.365	3.91%
1,2-Dichloropropane	0.416	0.373	0.376	0.400	0.405	0.409	0.396	4.49%
Methylmethacrylate	0.232	0.261	0.287	0.331	0.348	0.339	0.300	15.78% NT
Bromodichloromethane	0.461	0.427	0.453	0.477	0.494	0.506	0.470	6.12%
Dibromomethane	0.277	0.266	0.291	0.297	0.306	0.299	0.289	5.19%
1,4-Dioxane	0.003	0.004	0.005	0.006	0.006	0.006	0.005	19.13% NT
2-Chloroethylvinyl ether	0.003	0.004	0.004				0.004	8.23%
2-Hexanone	0.157	0.168	0.181	0.212	0.223	0.214	0.193	14.23%
Methyl isobutyl ketone	0.176	0.172	0.187	0.211	0.224	0.219	0.198	11.44%
cis-1,3-Dichloropropene	0.558	0.540	0.541	0.599	0.637	0.658	0.589	8.66%
d8-Toluene	1.285	1.266	1.264	1.286	1.298	1.284	1.281	1.01%
Toluene	0.906	0.783	0.805	0.838	0.876	0.903	0.852	6.03%
trans-1,3-Dichloropropene	0.483	0.460	0.480	0.553	0.612	0.643	0.539	14.17%
1,1,2-Trichloroethane	0.285	0.311	0.299	0.313	0.329	0.329	0.311	5.55%
1,3-Dichloropropane	0.607	0.585	0.592	0.627	0.649	0.654	0.619	4.67%
Tetrachloroethene	0.331	0.292	0.295	0.302	0.323	0.332	0.313	5.86%
Dibromochloromethane	0.263	0.250	0.269	0.305	0.337	0.351	0.296	14.09%

1,2-Dibromoethane	0.330	0.340	0.336	0.372	0.386	0.388	0.359	7.36%
Chlorobenzene	1.044	0.931	0.943	0.958	1.016	1.045	0.989	5.21%
1,1,1,2-Tetrachloroethane	0.258	0.258	0.261	0.296	0.328	0.349	0.292	13.60%
Ethylbenzene	1.892	1.688	1.668	1.762	1.867	1.934	1.802	6.20%
m,p-Xylene	0.628	0.567	0.581	0.594	0.653	0.702	0.621	8.18%
o-Xylene	0.592	0.535	0.546	0.576	0.611	0.637	0.583	6.66%
Styrene	1.014	0.959	0.988	1.077	1.157	1.192	1.064	8.86%
Bromoform	0.182	0.202	0.221	0.272	0.328	0.347	0.258	26.42%
Isopropylbenzene	1.328	1.228	1.266	1.338	1.428	1.506	1.349	7.63%
1,1,2,2-Tetrachloroethane	0.571	0.539	0.567	0.622	0.647	0.613	0.593	6.84%
Bromofluorobenzene	0.555	0.547	0.554	0.556	0.554	0.545	0.552	0.82%
1,2,3-Trichloropropane	0.439	0.458	0.448	0.473	0.491	0.472	0.464	4.09%
trans-1,4-Dichloro-2-butene	0.107	0.120	0.152	0.179	0.190		0.149	24.08% NT
n-Propylbenzene	1.812	1.721	1.717	1.803	1.955	2.067	1.846	7.50%
Bromobenzene	0.423	0.411	0.416	0.440	0.472	0.486	0.441	7.00%
1,3,5-Trimethylbenzene	1.048	0.999	1.030	1.082	1.164	1.247	1.095	8.55%
2-Chlorotoluene	1.245	1.211	1.203	1.267	1.382	1.423	1.289	7.14%
4-Chlorotoluene	1.226	1.093	1.118	1.165	1.194	1.278	1.179	5.82%
tert-butylbenzene	0.860	0.733	0.757	0.799	0.823	0.875	0.808	6.93%
1,2,4-Trimethylbenzene	1.171	1.053	1.040	1.123	1.217	1.293	1.149	8.51%
sec-butylbenzene	1.138	1.116	1.123	1.181	1.262	1.307	1.188	6.69%
p-isopropyltoluene	0.848	0.846	0.866	0.905	0.976	1.028	0.912	8.25%
1,3-Dichlorobenzene	0.695	0.693	0.692	0.725	0.774	0.806	0.731	6.67%
1,4-Dichlorobenzene	1.479	1.275	1.329	1.365	1.440	1.475	1.394	6.02%
n-butylbenzene	1.703	1.599	1.654	1.734	1.862	1.896	1.741	6.71%
1,2-Dichlorobenzene	1.273	1.239	1.227	1.266	1.364	1.392	1.294	5.27%
1,2-Dibromo-3-chloropropane	0.117	0.150	0.156	0.189	0.202	0.190	0.167	19.23%
1,2,4-Trichlorobenzene	0.571	0.616	0.606	0.649	0.706	0.741	0.648	9.90%
Hexachlorobutadiene	0.108	0.097	0.121	0.124	0.135	0.144	0.121	14.11%
Naphthalene	1.696	1.757	1.879	2.123	2.255	2.238	1.991	12.34%
1,2,3-Trichlorobenzene	0.574	0.552	0.589	0.627	0.685	0.705	0.622	9.99%
1,3,5-Trichlorobenzene	0.669	0.641	0.635	0.667	0.719	0.751	0.680	6.68%
							AVG	**10.09%**

Response Factor Report Instrumen

Method : C:\HPCHEM\1\METHODS\V805266B.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Tue May 30 08:22:32 2006
Response via : Initial Calibration

5-30-06 SL ✓

5-30-06

Calibration Files
1 =B36408SI.D 2 =B36409SI.D 5 =B36410SI.D
20 =B36411SI.D 50 =B36412SI.D 100 =B36413SI.D

		Compound	1	2	5	20	50	100	Avg	%RSD
1)	I	Pentafluorobenzene	ISTD							
2)	TM	Dichlorodifluoromet	0.695	0.687	0.694	0.688	0.676	0.656	0.683 ✓	2.17
3)	TMP	Chloromethane		1.382	1.290	1.129	1.164	1.097	1.212 ✓	9.88
4)	TMC	Vinyl Chloride		0.882	0.791	0.585	0.553	0.512	0.665	24.40# ✓
5)	TM	Bromomethane	0.343	0.262	0.320	0.447	0.583	0.591	0.424	32.86
6)	TM	Chloroethane	0.763	0.631	0.618	0.603	0.578	0.540	0.622	12.24
7)	TM	t-Butyl alcohol (TB		0.050	0.056	0.065	0.073	0.068	0.062#	14.99
8)	TM	Trichlorofluorometh	1.190	1.175	1.231	1.180	1.168	1.125	1.178	2.90
9)	TM	Diethyl ether	0.439	0.468	0.450	0.522	0.489	0.437	0.467	7.07
10)	TM	1,1,2-Trichlorotrif	0.504	0.499	0.518	0.516	0.516	0.510	0.511	1.52
11)	T	Acrolein	0.323	0.367	0.361	0.387	0.394	0.375	0.368	6.86
12)	TM	Acetone	0.331	0.297	0.299	0.307	0.304	0.281	0.303	5.34
13)	TMC	1,1-Dichloroethene	0.471	0.449	0.461	0.480	0.488	0.493	0.474	3.54# ✓
14)	TM	Methyl iodide	0.180	0.245	0.400	0.572	0.681	0.697	0.462	~~47.98~~ Q
15)	TM	Di-isopropyl ether	2.597	2.385	2.445	2.466	2.505	2.509	2.484	2.88
16)	TM	Methylene Chloride		0.749	0.674	0.630	0.629	0.621	0.660	8.10
17)	TM	Carbon Disulfide	1.904	1.779	1.798	1.869	1.939	1.960	1.875	3.95
18)	TM	Acrylonitrile	0.283	0.319	0.305	0.337	0.337	0.325	0.318	6.53
19)	TM	Methyl-tert-butyl e	1.763	1.701	1.750	1.786	1.778	1.754	1.755	1.70
20)	TM	trans-1,2-Dichloroe	0.606	0.539	0.525	0.548	0.549	0.555	0.553 ✓	5.04
21)	TMP	1,1-Dichloroethane	1.145	1.095	1.081	1.090	1.114	1.124	1.108 ✓	2.17
22)	TM	Vinyl acetate	0.030	0.060	0.079	0.104	0.117	0.116	0.084#	~~41.36~~ Q
23)	TM	Methyl ethyl ketone	0.429	0.436	0.447	0.486	0.474	0.443	0.453	4.92
24)	TM	Ethyl t-butyl ether	2.027	2.007	1.934	2.042	2.033	2.056	2.017	2.16
25)	TM	2,2-Dichloropropane	0.696	0.631	0.644	0.719	0.765	0.767	0.703	8.28
26)	TM	cis-1,2-Dichloroeth	0.593	0.566	0.566	0.586	0.598	0.599	0.585	2.60
27)	TM	t-Amyl methyl ether	1.688	1.637	1.655	1.727	1.733	1.724	1.694	2.40
28)	S	DB HOLDER #28	0.829	0.828	0.820	0.806	0.759	0.768	0.802	3.84
29)	TMC	Chloroform	1.188	1.031	1.055	1.065	1.066	1.067	1.079	5.13# ✓
30)	TM	Bromochloromethane	0.259	0.259	0.263	0.263	0.263	0.261	0.261	0.77
31)	TM	Tetrahydrofuran		0.092	0.084	0.092	0.092	0.086	0.089#	4.22
32)	TM	1,1,1-Trichloroetha	0.794	0.729	0.741	0.762	0.769	0.786	0.763	3.27
33)	T	NO COMPOUND							0.000#	-1.00
34)	TM	1,1-Dichloropropene	0.879	0.813	0.799	0.817	0.806	0.809	0.821	3.54
35)	TM	Carbon Tetrachlorid	0.465	0.432	0.479	0.507	0.525	0.545	0.492	8.46
36)	S	d4-1,2-Dichloroetha	0.829	0.828	0.820	0.806	0.759	0.768	0.802	3.84
37)	TM	1,2-Dichloroethane	1.129	1.043	1.005	0.979	0.966	0.949	1.012	6.54
38)	TM	Benzene	2.566	2.235	2.216	2.282	2.272	2.322	2.316	5.53
39)	I	1,4-Difluorobenzene	ISTD							
40)	TM	Trichloroethene	0.382	0.361	0.342	0.358	0.371	0.376	0.365	3.91 ✓
41)	TMC	1,2-Dichloropropane	0.416	0.373	0.376	0.400	0.405	0.409	0.396	4.49# ✓
42)	TM	Methylmethacrylate	0.232	0.261	0.287	0.331	0.348	0.339	0.300	15.78 NT

(#) = Out of Range

Method : C:\HPCHEM\1\METHODS\V805266B.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Tue May 30 08:22:32 2006
Response via : Initial Calibration

Calibration Files
1 =B36408SI.D 2 =B36409SI.D 5 =B36410SI.D
20 =B36411SI.D 50 =B36412SI.D 100 =B36413SI.D

		Compound	1	2	5	20	50	100	Avg	%RSD
43)	T	Methyl acetate	0.394	0.414	0.428	0.474	0.490	0.477	0.446	8.79
44)	T	Cyclohexane	0.525	0.517	0.525	0.551	0.559	0.575	0.542	4.29
45)	TM	Bromodichloromethan	0.461	0.427	0.453	0.477	0.494	0.506	0.470	6.12
46)	TM	Dibromomethane	0.277	0.266	0.291	0.297	0.306	0.299	0.289	5.19
47)	TM	1,4-Dioxane	0.003	0.004	0.005	0.006	0.006	0.006	0.005#	19.13 NT
48)	T	Methylcyclohexane	0.350	0.358	0.369	0.390	0.402	0.407	0.379	6.22
49)	T	NO COMPOUND							0.000#	-1.00
50)	T	2-Chloroethylvinyl				0.003	0.004	0.004	0.004#	8.23
51)	TM	2-Hexanone	0.157	0.168	0.181	0.212	0.223	0.214	0.193	14.23
52)	TM	Methyl isobutyl ket	0.176	0.172	0.187	0.211	0.224	0.219	0.198	11.44
53)	TM	cis-1,3-Dichloropro	0.558	0.540	0.541	0.599	0.637	0.658	0.589	8.66
54)	S	d8-Toluene	1.285	1.266	1.264	1.286	1.298	1.284	1.281	1.01
55)	TMC	Toluene	0.906	0.783	0.805	0.838	0.876	0.903	0.852	6.03#
56)	TM	trans-1,3-Dichlorop	0.483	0.460	0.480	0.553	0.612	0.643	0.539	14.17
57)	T	NO COMPOUND							0.000#	-1.00
58)	TM	1,1,2-Trichloroetha	0.285	0.311	0.299	0.313	0.329	0.329	0.311	5.55
59)	TM	1,3-Dichloropropane	0.607	0.585	0.592	0.627	0.649	0.654	0.619	4.67
60)	TM	Tetrachloroethene	0.331	0.292	0.295	0.302	0.323	0.332	0.313	5.86
61)	TM	Dibromochloromethan	0.263	0.250	0.269	0.305	0.337	0.351	0.296	14.09
62)	TM	1,2-Dibromoethane	0.330	0.340	0.336	0.372	0.386	0.388	0.359	7.36
63)	I	*d5-Chlorobenzene	----------ISTD----------							
64)	TMP	Chlorobenzene	1.044	0.931	0.943	0.958	1.016	1.045	0.989	5.21
65)	TM	1,1,1,2-Tetrachloro	0.258	0.258	0.261	0.296	0.328	0.349	0.292	13.60
66)	TMC	Ethylbenzene	1.892	1.688	1.668	1.762	1.867	1.934	1.802	6.20#
67)	TM	m,p-Xylene	0.628	0.567	0.581	0.594	0.653	0.702	0.621	8.18
68)	TM	o-Xylene	0.592	0.535	0.546	0.576	0.611	0.637	0.583	6.66
69)	TM	Styrene	1.014	0.959	0.988	1.077	1.157	1.192	1.064	8.86
70)	TMP	Bromoform	0.182	0.202	0.221	0.272	0.328	0.347	0.258	26.42
71)	TM	Isopropylbenzene	1.328	1.228	1.266	1.338	1.428	1.506	1.349	7.63
72)	TMP	1,1,2,2-Tetrachloro	0.571	0.539	0.567	0.622	0.647	0.613	0.593	6.84
73)	S	Bromofluorobenzene	0.555	0.547	0.554	0.556	0.554	0.545	0.552	0.82
74)	TM	1,2,3-Trichloroprop	0.439	0.458	0.448	0.473	0.491	0.472	0.464	4.09
75)	TM	trans-1,4-Dichloro-		0.107	0.120	0.152	0.179	0.190	0.149	24.08
76)	TM	n-Propylbenzene	1.812	1.721	1.717	1.803	1.955	2.067	1.846	7.50
77)	TM	Bromobenzene	0.423	0.411	0.416	0.440	0.472	0.486	0.441	7.00
78)	TM	1,3,5-Trimethylbenz	1.048	0.999	1.030	1.082	1.164	1.247	1.095	8.55
79)	TM	2-Chlorotoluene	1.245	1.211	1.203	1.267	1.382	1.423	1.289	7.14
80)	TM	4-Chlorotoluene	1.226	1.093	1.118	1.165	1.194	1.278	1.179	5.82
81)	TM	tert-butylbenzene	0.860	0.733	0.757	0.799	0.823	0.875	0.808	6.93
82)	TM	1,2,4-Trimethylbenz	1.171	1.053	1.040	1.123	1.217	1.293	1.149	8.51
83)	T	Diisobutylketone	1.111	1.124	1.164	1.299	1.469	1.488	1.276	13.36
84)	TM	sec-butylbenzene	1.138	1.116	1.123	1.181	1.262	1.307	1.188	6.69
85)	TM	p-isopropyltoluene	0.848	0.846	0.866	0.905	0.976	1.028	0.912	8.25

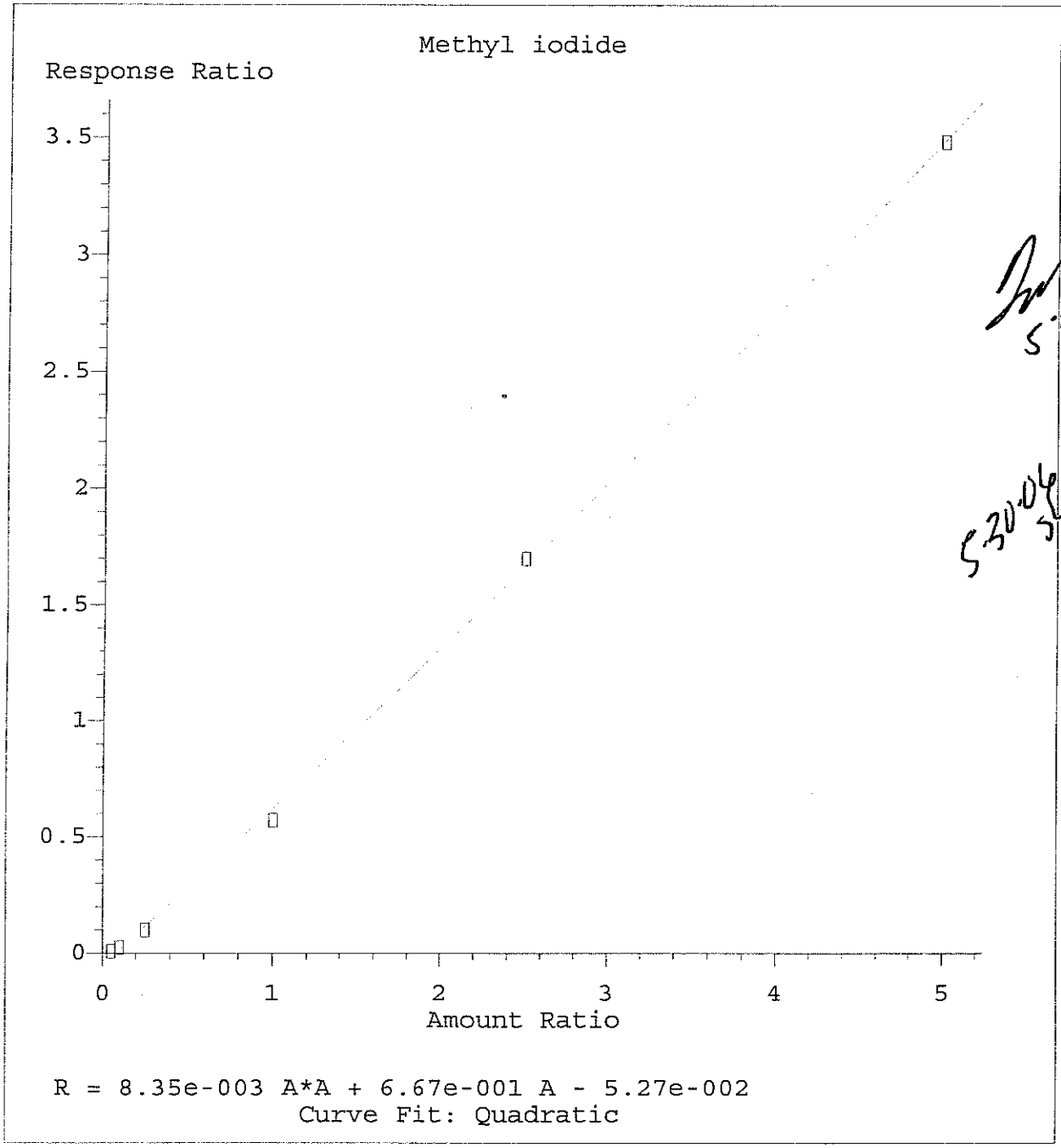
(#) = Out of Range

Method : C:\HPCHEM\1\METHODS\V805266B.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Tue May 30 08:22:32 2006
Response via : Initial Calibration

Calibration Files
1 =B36408SI.D 2 =B36409SI.D 5 =B36410SI.D
20 =B36411SI.D 50 =B36412SI.D 100 =B36413SI.D

		Compound	1	2	5	20	50	100	Avg	%RSD
86)	TM	1,3-Dichlorobenzene	0.695	0.693	0.692	0.725	0.774	0.806	0.731	6.67
87)	I	1,4-Dichlorobenzene-d	----------	----------	----------	ISTD	----------	----------		
88)	TM	1,4-Dichlorobenzene	1.479	1.275	1.329	1.365	1.440	1.475	1.394	6.02
89)	TM	n-butylbenzene	1.703	1.599	1.654	1.734	1.862	1.896	1.741	6.71
90)	TM	1,2-Dichlorobenzene	1.273	1.239	1.227	1.266	1.364	1.392	1.294	5.27
91)	T	NO COMPOUND							0.000#	-1.00
92)	TM	1,2-Dibromo-3-chlor	0.117	0.150	0.156	0.189	0.202	0.190	0.167	19.23
93)	T	NO COMPOUND							0.000#	-1.00
94)	TM	1,2,4-Trichlorobenz	0.571	0.616	0.606	0.649	0.706	0.741	0.648	9.90
95)	TM	Hexachlorobutadiene	0.108	0.097	0.121	0.124	0.135	0.144	0.121	14.11
96)	TM	Naphthalene	1.696	1.757	1.879	2.123	2.255	2.238	1.991	12.34
97)	TM	1,2,3-Trichlorobenz	0.574	0.552	0.589	0.627	0.685	0.705	0.622	9.99
98)	TM	1,3,5-Trichlorobenz	0.669	0.641	0.635	0.667	0.719	0.751	0.680	6.68

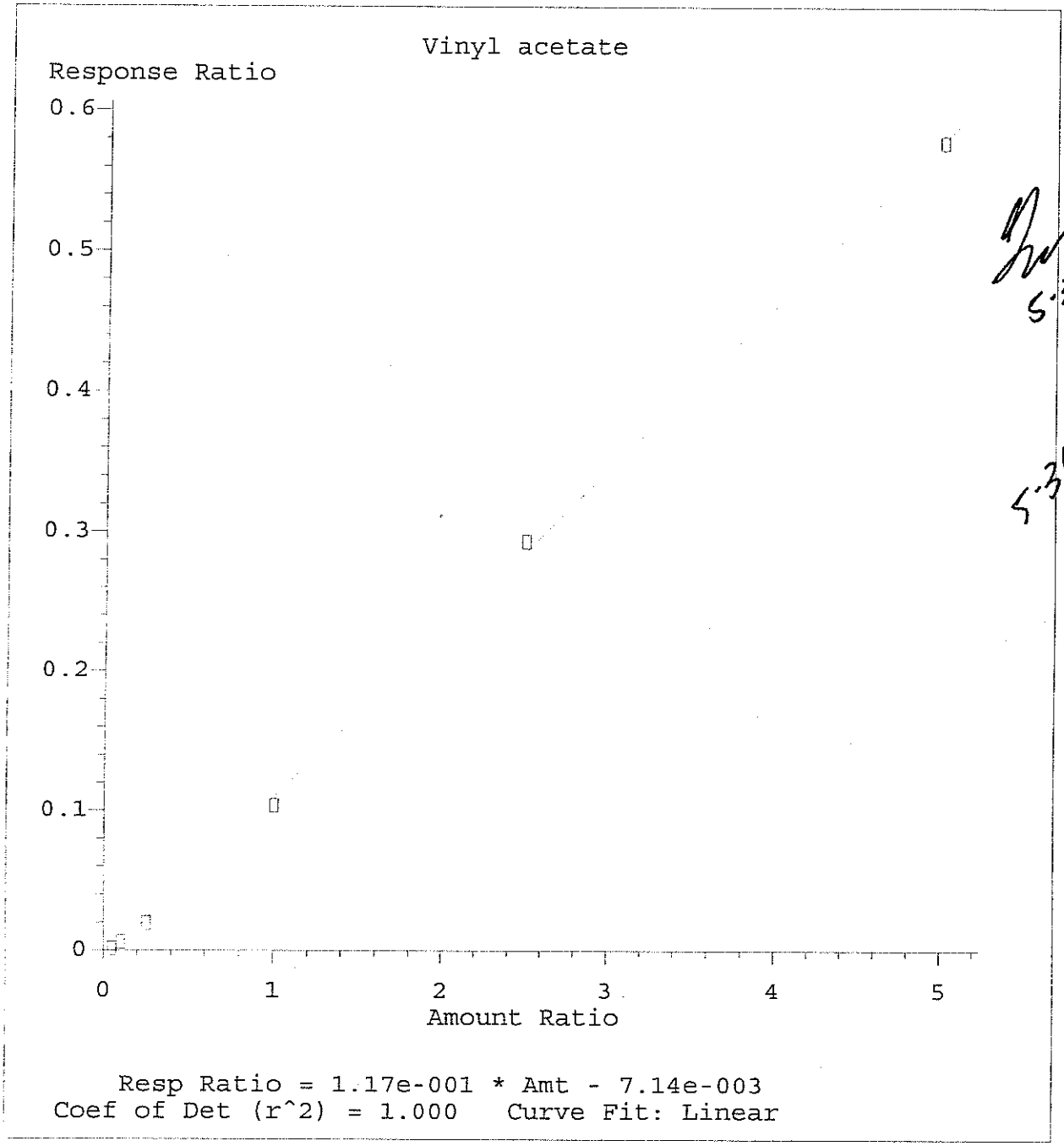
(#) = Out of Range



5-30-06

5.30.06

Method Name: C:\HPCHEM\1\METHODS\V805266B.M
Calibration Table Last Updated: Tue May 30 08:15:26 2006



5-30-06

5-30-06

Method Name: C:\HPCHEM\1\METHODS\V805266B.M
Calibration Table Last Updated: Tue May 30 08:15:26 2006

Quantitation Report (Not Reviewed)

Data File : C:\HPCHEM\1\DATA\052606-B\B36408SI.D Vial: 2
Acq On : 26 May 2006 10:18 am Operator:
Sample : 1 PPB Inst : Instrumen
Misc : 5000 Multiplr: 1.00
Integration Params: rteint.p
Quant Time: May 30 8:17 2006 Quant Results File: V805266B.RES

Quant Method : C:\HPCHEM\1\METHODS\V805266B.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Tue May 30 08:15:26 2006
Response via : Initial Calibration
DataAcq Meth : V805266B

5-30-06 SV 5-30-06

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.11	168	92853	20.00	ug/L	0.00
39) 1,4-Difluorobenzene	8.54	114	147055	20.00	ug/L	0.00
63) *d5-Chlorobenzene	12.84	117	124955	20.00	ug/L	0.00
87) 1,4-Dichlorobenzene-d4	16.33	152	68277	20.00	ug/L	0.00
System Monitoring Compounds						
28) DB HOLDER #28	7.96	65	76955	20.68	ug/L	0.00
Spiked Amount 20.000	Range 70 - 130		Recovery = 103.40%			
36) d4-1,2-Dichloroethane	7.96	65	76955	20.68	ug/L	0.00
Spiked Amount 20.000	Range 70 - 130		Recovery = 103.40%			
54) d8-Toluene	10.69	98	188913	20.06	ug/L	0.00
Spiked Amount 20.000	Range 70 - 130		Recovery = 100.30%			
73) Bromofluorobenzene	14.60	95	69374	20.12	ug/L	0.00
Spiked Amount 20.000	Range 70 - 130		Recovery = 100.60%			
Target Compounds						Qvalue
2) Dichlorodifluoromethane	1.70	85	3227	1.02	ug/L	98
3) Chloromethane	1.97	50	8179	1.45	ug/L	84
4) Vinyl Chloride	2.05	62	4856	1.57	ug/L	98
5) Bromomethane	2.55	94	1592	0.81	ug/L	87
6) Chloroethane	2.65	64	3542	1.23	ug/L	99
8) Trichlorofluoromethane	2.96	101	5526	1.01	ug/L	96
10) 1,1,2-Trichlorotrifluoroet	3.57	101	2341	0.99	ug/L	92
12) Acetone	3.65	43	7676	5.45	ug/L	95
13) 1,1-Dichloroethene	3.83	96	2186	0.99	ug/L #	80
14) Methyl iodide	4.28	142	834	1.85	ug/L #	60
15) Di-isopropyl ether (DIPE)	5.62	45	12058	1.05	ug/L	95
17) Carbon Disulfide	4.57	76	8838	1.02	ug/L	98
19) Methyl-tert-butyl ether (M	4.88	73	8186	1.00	ug/L	96
20) trans-1,2-Dichloroethene	5.07	96	2814	1.10	ug/L	93
21) 1,1-Dichloroethane	5.75	63	5316	1.03	ug/L	88
22) Vinyl acetate	5.80	86	138	1.47	ug/L #	1
24) Ethyl t-butyl ether (ETBE)	6.26	59	9409	1.01	ug/L	97
25) 2,2-Dichloropropane	6.61	77	3232	0.99	ug/L	90
26) cis-1,2-Dichloroethene	6.68	96	2752	1.01	ug/L	93
27) t-Amyl methyl ether (TAME)	7.93	73	7836	1.00	ug/L	97
29) Chloroform	6.91	83	5515	1.10	ug/L	90
30) Bromochloromethane	7.14	128	1204	0.99	ug/L	93
32) 1,1,1-Trichloroethane	7.49	97	3684	1.04	ug/L	94
34) 1,1-Dichloropropene	7.72	75	4079	1.07	ug/L	93
35) Carbon Tetrachloride	7.86	119	2160	0.95	ug/L	88
37) 1,2-Dichloroethane	8.09	62	5242	1.12	ug/L	98

(#) = qualifier out of range (m) = manual integration

Quantitation Report (Not Reviewed)

Data File : C:\HPCHEM\1\DATA\052606-B\B36408SI.D Vial: 2
Acq On : 26 May 2006 10:18 am Operator:
Sample : 1 PPB Inst : Instrumen
Misc : 5000 Multiplr: 1.00
Integration Params: rteint.p
Quant Time: May 30 8:17 2006 Quant Results File: V805266B.RES

Quant Method : C:\HPCHEM\1\METHODS\V805266B.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Tue May 30 08:15:26 2006
Response via : Initial Calibration
DataAcq Meth : V805266B

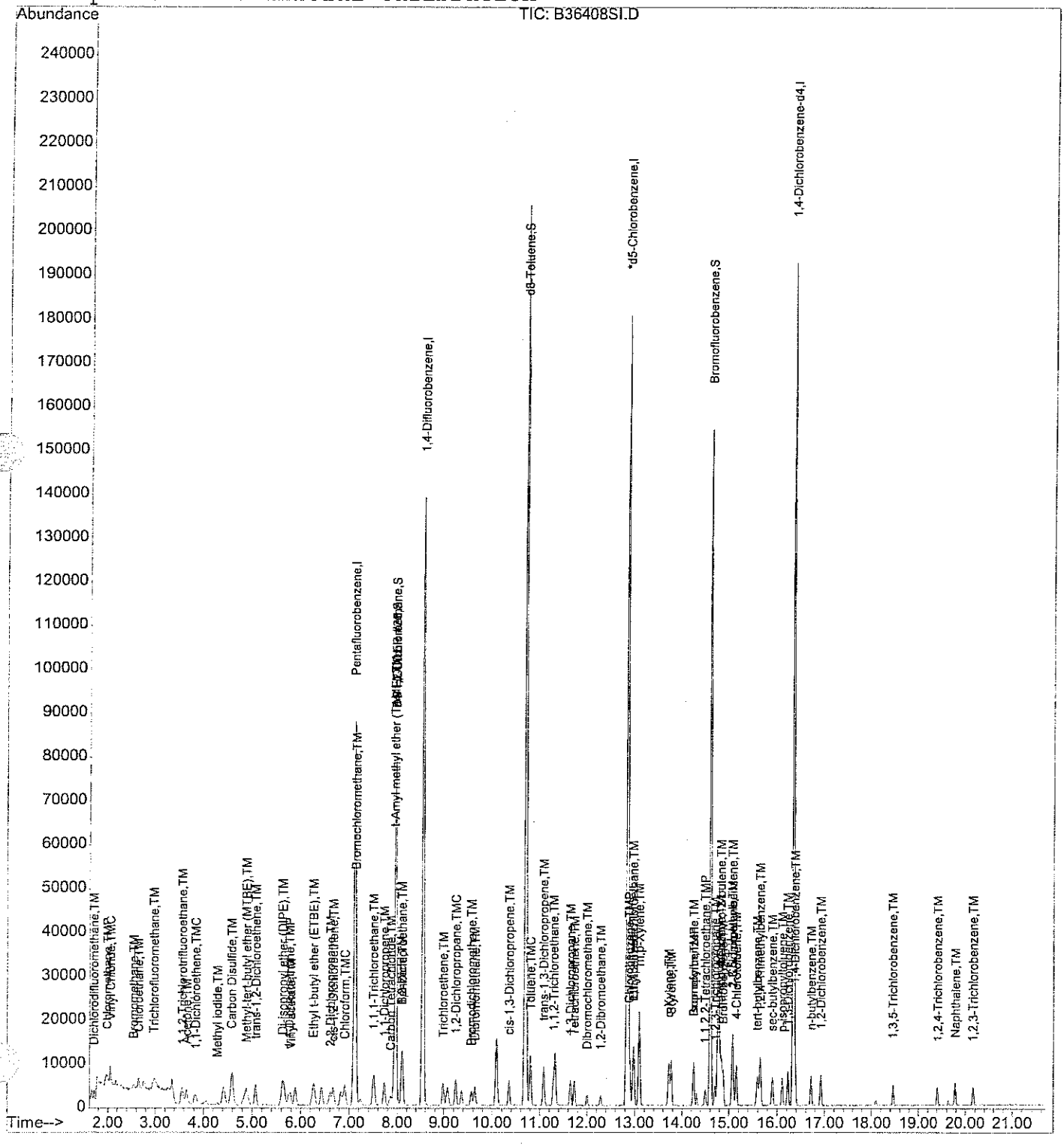
	Compound	R.T.	QIon	Response	Conc Unit	Qvalue
38)	Benzene	8.11	78	11912	1.11 ug/L	99
40)	Trichloroethene	9.00	95	2806	1.05 ug/L	94
41)	1,2-Dichloropropane	9.24	63	3060	1.05 ug/L	90
45)	Bromodichloromethane	9.57	83	3390	0.98 ug/L	89
46)	Dibromomethane	9.64	93	2038	0.96 ug/L	92
53)	cis-1,3-Dichloropropene	10.36	75	4103	0.95 ug/L	99
55)	Toluene	10.80	92	6658	1.06 ug/L	88
56)	trans-1,3-Dichloropropene	11.07	75	3548	0.90 ug/L	92
58)	1,1,2-Trichloroethane	11.28	83	2095	0.92 ug/L	90
59)	1,3-Dichloropropane	11.65	76	4465	0.98 ug/L	94
60)	Tetrachloroethene	11.73	166	2431	1.06 ug/L	93
61)	Dibromochloromethane	12.00	129	1935	0.89 ug/L	91
62)	1,2-Dibromoethane	12.29	107	2424	0.92 ug/L	87
64)	Chlorobenzene	12.89	112	6520	1.05 ug/L	93
65)	1,1,1,2-Tetrachloroethane	12.96	131	1613	0.89 ug/L	97
66)	Ethylbenzene	12.98	91	11820	1.05 ug/L	99
67)	m,p-Xylene	13.10	106	7849	2.02 ug/L	96
68)	o-Xylene	13.72	106	3701	1.02 ug/L #	85
69)	Styrene	13.77	104	6337	0.95 ug/L	99
70)	Bromoform	14.23	173	1135	0.70 ug/L #	27
71)	Isopropylbenzene	14.24	105	8294	0.98 ug/L	92
72)	1,1,2,2-Tetrachloroethane	14.49	83	3566	0.96 ug/L	96
74)	1,2,3-Trichloropropane	14.71	75	2745	0.95 ug/L #	84
75)	trans-1,4-Dichloro-2-buten	14.83	89	627	0.67 ug/L #	35
76)	n-Propylbenzene	14.83	91	11324	0.98 ug/L	93
77)	Bromobenzene	14.87	156	2641	0.96 ug/L	91
78)	1,3,5-Trimethylbenzene	15.06	105	6545	0.96 ug/L	96
79)	2-Chlorotoluene	15.07	91	7777	0.97 ug/L	99
80)	4-Chlorotoluene	15.15	91	7658	1.04 ug/L	96
81)	tert-butylbenzene	15.58	119	5370	1.06 ug/L	99
82)	1,2,4-Trimethylbenzene	15.64	105	7316	1.02 ug/L	94
84)	sec-butylbenzene	15.90	105	7107	0.96 ug/L	98
85)	p-isopropyltoluene	16.11	119	5296	0.93 ug/L	95
86)	1,3-Dichlorobenzene	16.22	146	4344	0.95 ug/L	98
88)	1,4-Dichlorobenzene	16.38	146	5050	1.06 ug/L	94
89)	n-butylbenzene	16.72	91	5813	0.98 ug/L	99
90)	1,2-Dichlorobenzene	16.92	146	4347	0.98 ug/L	85
94)	1,2,4-Trichlorobenzene	19.41	180	1951	0.88 ug/L	94
96)	Naphthalene	19.77	128	5791	0.85 ug/L	100
97)	1,2,3-Trichlorobenzene	20.16	180	1958	0.92 ug/L	91
98)	1,3,5-Trichlorobenzene	18.45	180	2283	0.98 ug/L	93

(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : C:\HPCHEM\1\DATA\052606-B\B36408SI.D Vial: 2
Acq On : 26 May 2006 10:18 am Operator:
Sample : 1 PPB Inst : Instrumen
Misc : 5000 Multiplr: 1.00
MS Integration Params: rteint.p
Quant Time: May 30 8:17 2006 Quant Results File: V805266B.RES

Method : C:\HPCHEM\1\METHODS\V805266B.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Tue May 30 08:15:26 2006
Response via : Initial Calibration



Quantitation Report (Not Reviewed)

Data File : C:\HPCHEM\1\DATA\052606-B\B36409SI.D Vial: 3
Acq On : 26 May 2006 10:53 am Operator:
Sample : 2 PPB Inst : Instrumen
Misc : 5000 Multiplr: 1.00
Integration Params: rteint.p
Quant Time: May 30 8:17 2006 Quant Results File: V805266B.RES

Quant Method : C:\HPCHEM\1\METHODS\V805266B.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Tue May 30 08:15:26 2006
Response via : Initial Calibration
DataAcq Meth : V805266B

5-30-06 SL 5-30-06

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.11	168	93127	20.00	ug/L	0.00
39) 1,4-Difluorobenzene	8.54	114	147758	20.00	ug/L	0.00
63) *d5-Chlorobenzene	12.84	117	124425	20.00	ug/L	0.00
87) 1,4-Dichlorobenzene-d4	16.34	152	69352	20.00	ug/L	0.00
System Monitoring Compounds						
28) DB HOLDER #28	7.96	65	77115	20.66	ug/L	0.00
Spiked Amount 20.000	Range 70 - 130		Recovery = 103.30%			
36) d4-1,2-Dichloroethane	7.96	65	77115	20.66	ug/L	0.00
Spiked Amount 20.000	Range 70 - 130		Recovery = 103.30%			
54) d8-Toluene	10.69	98	187117	19.78	ug/L	0.00
Spiked Amount 20.000	Range 70 - 130		Recovery = 98.90%			
73) Bromofluorobenzene	14.60	95	68031	19.82	ug/L	0.00
Spiked Amount 20.000	Range 70 - 130		Recovery = 99.10%			
Target Compounds						Qvalue
2) Dichlorodifluoromethane	1.70	85	6395	2.01	ug/L	96
3) Chloromethane	1.98	50	12872	2.28	ug/L	96
4) Vinyl Chloride	2.06	62	8212	2.65	ug/L	94
5) Bromomethane	2.55	94	2441	1.24	ug/L	95
6) Chloroethane	2.64	64	5875	2.03	ug/L	97
8) Trichlorofluoromethane	2.98	101	10946	1.99	ug/L	89
9) Diethyl ether	3.36	74	4355	2.00	ug/L #	88
10) 1,1,2-Trichlorotrifluoroet	3.56	101	4644	1.95	ug/L	98
12) Acetone	3.66	43	13812	9.79	ug/L	99
13) 1,1-Dichloroethene	3.82	96	4181	1.90	ug/L	96
14) Methyl iodide	4.29	142	2277	2.31	ug/L #	90
15) Di-isopropyl ether (DIPE)	5.63	45	22209	1.92	ug/L	100
16) Methylene Chloride	4.59	84	6972	2.27	ug/L	95
17) Carbon Disulfide	4.56	76	16568	1.90	ug/L #	94
18) Acrylonitrile	4.81	53	2969	2.01	ug/L	97
19) Methyl-tert-butyl ether (M	4.88	73	15841	1.94	ug/L	98
20) trans-1,2-Dichloroethene	5.08	96	5016	1.95	ug/L	89
21) 1,1-Dichloroethane	5.76	63	10199	1.98	ug/L	92
22) Vinyl acetate	5.80	86	556	2.23	ug/L #	17
23) Methyl ethyl ketone	6.43	43	20318	9.64	ug/L	99
24) Ethyl t-butyl ether (ETBE)	6.27	59	18694	1.99	ug/L	99
25) 2,2-Dichloropropane	6.60	77	5872	1.79	ug/L	94
26) cis-1,2-Dichloroethene	6.67	96	5274	1.94	ug/L	93
27) t-Amyl methyl ether (TAME)	7.93	73	15247	1.93	ug/L	98
29) Chloroform	6.91	83	9599	1.91	ug/L	92
30) Bromochloromethane	7.15	128	2409	1.98	ug/L #	84

(#) = qualifier out of range (m) = manual integration

Quantitation Report (Not Reviewed)

Data File : C:\HPCHEM\1\DATA\052606-B\B36409SI.D Vial: 3
Acq On : 26 May 2006 10:53 am Operator:
Sample : 2 PPB Inst : Instrumen
Misc : 5000 Multiplr: 1.00
Integration Params: rteint.p
Quant Time: May 30 8:17 2006 Quant Results File: V805266B.RES

Quant Method : C:\HPCHEM\1\METHODS\V805266B.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Tue May 30 08:15:26 2006
Response via : Initial Calibration
DataAcq Meth : V805266B

	Compound	R.T.	QIon	Response	Conc Unit	Qvalue
32)	1,1,1-Trichloroethane	7.49	97	6790	1.91 ug/L	95
34)	1,1-Dichloropropene	7.73	75	7567	1.98 ug/L	95
35)	Carbon Tetrachloride	7.86	119	4021	1.75 ug/L	93
37)	1,2-Dichloroethane	8.09	62	9714	2.06 ug/L	97
38)	Benzene	8.10	78	20818	1.93 ug/L	96
40)	Trichloroethene	8.99	95	5334	1.98 ug/L	95
41)	1,2-Dichloropropane	9.25	63	5516	1.88 ug/L	94
42)	Methylmethacrylate	9.37	69	3861	1.74 ug/L	93
43)	Methyl acetate	4.40	43	6120	1.86 ug/L #	94
44)	Cyclohexane	7.51	56	7635	1.91 ug/L	99
45)	Bromodichloromethane	9.57	83	6311	1.82 ug/L	99
46)	Dibromomethane	9.63	93	3927	1.84 ug/L	98
47)	1,4-Dioxane	9.62	88	1650	43.77 ug/L	95
48)	Methylcyclohexane	9.08	83	5294	1.89 ug/L	96
51)	2-Hexanone	11.32	58	12402	8.72 ug/L	93
52)	Methyl isobutyl ketone	10.09	58	12676	8.66 ug/L	100
53)	cis-1,3-Dichloropropene	10.35	75	7972	1.83 ug/L	97
55)	Toluene	10.81	92	11569	1.84 ug/L	96
56)	trans-1,3-Dichloropropene	11.07	75	6802	1.71 ug/L	97
58)	1,1,2-Trichloroethane	11.28	83	4599	2.00 ug/L	96
59)	1,3-Dichloropropane	11.64	76	8642	1.89 ug/L	99
60)	Tetrachloroethene	11.73	166	4320	1.87 ug/L	90
61)	Dibromochloromethane	12.01	129	3700	1.69 ug/L	93
62)	1,2-Dibromoethane	12.29	107	5019	1.89 ug/L	100
64)	Chlorobenzene	12.90	112	11586	1.88 ug/L	93
65)	1,1,1,2-Tetrachloroethane	12.96	131	3205	1.77 ug/L	90
66)	Ethylbenzene	12.99	91	21008	1.87 ug/L	98
67)	m,p-Xylene	13.09	106	14106	3.65 ug/L	93
68)	o-Xylene	13.72	106	6656	1.84 ug/L	98
69)	Styrene	13.77	104	11932	1.80 ug/L	100
70)	Bromoform	14.24	173	2511	1.56 ug/L	94
71)	Isopropylbenzene	14.25	105	15276	1.82 ug/L	100
72)	1,1,2,2-Tetrachloroethane	14.49	83	6702	1.82 ug/L	96
74)	1,2,3-Trichloropropane	14.70	75	5702	1.98 ug/L	96
75)	trans-1,4-Dichloro-2-buten	14.82	89	1328	1.43 ug/L	80
76)	n-Propylbenzene	14.83	91	21415	1.86 ug/L	97
77)	Bromobenzene	14.87	156	5119	1.86 ug/L #	87
78)	1,3,5-Trimethylbenzene	15.07	105	12424	1.82 ug/L	97
79)	2-Chlorotoluene	15.08	91	15071	1.88 ug/L	99
80)	4-Chlorotoluene	15.15	91	13597	1.85 ug/L	98
81)	tert-butylbenzene	15.58	119	9118	1.81 ug/L	98

(#) = qualifier out of range (m) = manual integration

Quantitation Report (Not Reviewed)

Data File : C:\HPCHEM\1\DATA\052606-B\B36409SI.D Vial: 3
Acq On : 26 May 2006 10:53 am Operator:
Sample : 2 PPB Inst : Instrumen
Misc : 5000 Multiplr: 1.00
, Integration Params: rteint.p
Quant Time: May 30 8:17 2006 Quant Results File: V805266B.RES

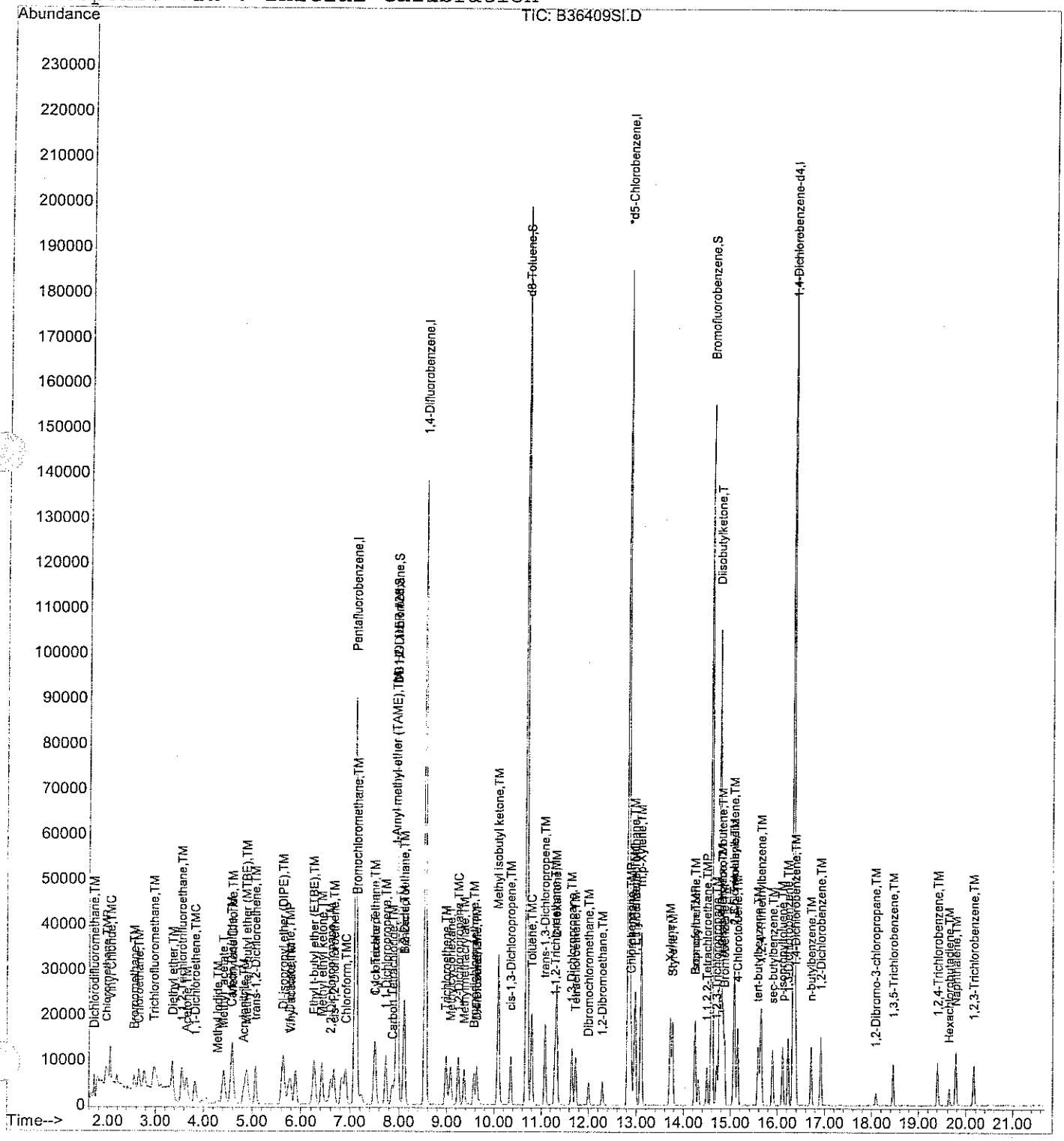
Quant Method : C:\HPCHEM\1\METHODS\V805266B.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Tue May 30 08:15:26 2006
Response via : Initial Calibration
DataAcq Meth : V805266B

	Compound	R.T.	QIon	Response	Conc	Unit		Qvalue
82)	1,2,4-Trimethylbenzene	15.64	105	13098	1.83	ug/L		97
83)	Diisobutylketone	14.77	57	69931	8.81	ug/L		96
84)	sec-butylbenzene	15.90	105	13889	1.88	ug/L		97
85)	p-isopropyltoluene	16.11	119	10532	1.86	ug/L		99
86)	1,3-Dichlorobenzene	16.23	146	8618	1.90	ug/L		98
88)	1,4-Dichlorobenzene	16.39	146	8843	1.83	ug/L		90
89)	n-butylbenzene	16.71	91	11086	1.84	ug/L		95
90)	1,2-Dichlorobenzene	16.92	146	8592	1.92	ug/L		92
92)	1,2-Dibromo-3-chloropropan	18.08	75	1037	1.79	ug/L	#	86
94)	1,2,4-Trichlorobenzene	19.41	180	4270	1.90	ug/L		89
95)	Hexachlorobutadiene	19.63	225	670	1.59	ug/L	#	46
96)	Naphthalene	19.78	128	12188	1.76	ug/L		100
97)	1,2,3-Trichlorobenzene	20.15	180	3827	1.77	ug/L		90
98)	1,3,5-Trichlorobenzene	18.46	180	4446	1.88	ug/L		91

(#) = qualifier out of range (m) = manual integration

Data File : C:\HPCHEM\1\DATA\052606-B\B36409SI.D Vial: 3
Acq On : 26 May 2006 10:53 am Operator:
Sample : 2 PPB Inst : Instrumen
Misc : 5000 Multiplr: 1.00
MS Integration Params: rteint.p
Quant Time: May 30 8:17 2006 Quant Results File: V805266B.RES

Method : C:\HPCHEM\1\METHODS\V805266B.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Tue May 30 08:15:26 2006
Response via : Initial Calibration



Quantitation Report (Not Reviewed)

Data File : C:\HPCHEM\1\DATA\052606-B\B36410SI.D Vial: 4
Acq On : 26 May 2006 11:22 am Operator:
Sample : 5 PPB Inst : Instrumen
Misc : 5000 Multiplr: 1.00
Integration Params: rteint.p
Quant Time: May 30 8:17 2006 Quant Results File: V805266B.RES

Quant Method : C:\HPCHEM\1\METHODS\V805266B.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Tue May 30 08:15:26 2006
Response via : Initial Calibration
DataAcq Meth : V805266B

5-30-06 SL ✓ 5-30-06

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.11	168	93722	20.00	ug/L	0.00
39) 1,4-Difluorobenzene	8.55	114	148737	20.00	ug/L	0.00
63) *d5-Chlorobenzene	12.84	117	126711	20.00	ug/L	0.00
87) 1,4-Dichlorobenzene-d4	16.34	152	70309	20.00	ug/L	0.00
System Monitoring Compounds						
28) DB HOLDER #28	7.96	65	76840	20.46	ug/L	0.00
Spiked Amount 20.000	Range 70 - 130		Recovery = 102.30%			
36) d4-1,2-Dichloroethane	7.96	65	76840	20.46	ug/L	0.00
Spiked Amount 20.000	Range 70 - 130		Recovery = 102.30%			
54) d8-Toluene	10.70	98	188046	19.74	ug/L	0.00
Spiked Amount 20.000	Range 70 - 130		Recovery = 98.70%			
73) Bromofluorobenzene	14.60	95	70194	20.08	ug/L	0.00
Spiked Amount 20.000	Range 70 - 130		Recovery = 100.40%			
Target Compounds						Qvalue
2) Dichlorodifluoromethane	1.70	85	16261	5.08	ug/L	95
3) Chloromethane	1.97	50	30218	5.32	ug/L	100
4) Vinyl Chloride	2.06	62	18539	5.95	ug/L	99
5) Bromomethane	2.55	94	7502	3.77	ug/L	93
6) Chloroethane	2.65	64	14474	4.96	ug/L	90
7) t-Butyl alcohol (TBA)	3.98	59	6540	22.42	ug/L #	86
8) Trichlorofluoromethane	2.96	101	28844	5.22	ug/L	98
9) Diethyl ether	3.37	74	10537	4.81	ug/L #	87
10) 1,1,2-Trichlorotrifluoroet	3.57	101	12140	5.07	ug/L	98
12) Acetone	3.65	43	35005	24.64	ug/L	99
13) 1,1-Dichloroethene	3.82	96	10802	4.87	ug/L	97
14) Methyl iodide	4.29	142	9378	4.56	ug/L	98
15) Di-isopropyl ether (DIPE)	5.64	45	57276	4.92	ug/L	99
16) Methylene Chloride	4.59	84	15781	5.10	ug/L	97
17) Carbon Disulfide	4.56	76	42126	4.79	ug/L	99
18) Acrylonitrile	4.82	53	7145	4.80	ug/L	97
19) Methyl-tert-butyl ether (M	4.88	73	41005	4.99	ug/L	99
20) trans-1,2-Dichloroethene	5.07	96	12290	4.74	ug/L	95
21) 1,1-Dichloroethane	5.75	63	25325	4.88	ug/L	99
22) Vinyl acetate	5.79	86	1846	4.57	ug/L	88
23) Methyl ethyl ketone	6.43	43	52415	24.71	ug/L	99
24) Ethyl t-butyl ether (ETBE)	6.27	59	45318	4.80	ug/L	98
25) 2,2-Dichloropropane	6.60	77	15078	4.57	ug/L	95
26) cis-1,2-Dichloroethene	6.67	96	13253	4.84	ug/L	95
27) t-Amyl methyl ether (TAME)	7.93	73	38788	4.89	ug/L	99
29) Chloroform	6.91	83	24719	4.89	ug/L	99

(#) = qualifier out of range (m) = manual integration

Quantitation Report (Not Reviewed)

Data File : C:\HPCHEM\1\DATA\052606-B\B36410SI.D Vial: 4
Acq On : 26 May 2006 11:22 am Operator:
Sample : 5 PPB Inst : Instrumen
Misc : 5000 Multiplr: 1.00
Integration Params: rteint.p
Quant Time: May 30 8:17 2006 Quant Results File: V805266B.RES

Quant Method : C:\HPCHEM\1\METHODS\V805266B.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Tue May 30 08:15:26 2006
Response via : Initial Calibration
DataAcq Meth : V805266B

Compound	R.T.	QIon	Response	Conc Unit	Qvalue
30) Bromochloromethane	7.15	128	6172	5.04 ug/L #	87
31) Tetrahydrofuran	7.22	71	1965	4.71 ug/L	94
32) 1,1,1-Trichloroethane	7.50	97	17364	4.85 ug/L	93
34) 1,1-Dichloropropene	7.74	75	18729	4.87 ug/L	97
35) Carbon Tetrachloride	7.87	119	11228	4.87 ug/L	98
37) 1,2-Dichloroethane	8.09	62	23540	4.97 ug/L	97
38) Benzene	8.10	78	51931	4.79 ug/L	99
40) Trichloroethene	8.99	95	12723	4.69 ug/L	95
41) 1,2-Dichloropropane	9.25	63	13978	4.74 ug/L	100
42) Methylmethacrylate	9.36	69	10665	4.78 ug/L	97
43) Methyl acetate	4.39	43	15920	4.80 ug/L	98
44) Cyclohexane	7.51	56	19527	4.84 ug/L	98
45) Bromodichloromethane	9.57	83	16848	4.82 ug/L	96
46) Dibromomethane	9.63	93	10828	5.03 ug/L	92
47) 1,4-Dioxane	9.62	88	4759	125.41 ug/L	94
48) Methylcyclohexane	9.08	83	13735	4.87 ug/L	98
51) 2-Hexanone	11.32	58	33717	23.55 ug/L	97
52) Methyl isobutyl ketone	10.08	58	34758	23.60 ug/L	99
53) cis-1,3-Dichloropropene	10.36	75	20105	4.59 ug/L	99
55) Toluene	10.80	92	29941	4.73 ug/L	98
56) trans-1,3-Dichloropropene	11.07	75	17861	4.46 ug/L	98
58) 1,1,2-Trichloroethane	11.29	83	11113	4.80 ug/L	97
59) 1,3-Dichloropropane	11.64	76	22028	4.78 ug/L	95
60) Tetrachloroethene	11.73	166	10951	4.71 ug/L	91
61) Dibromochloromethane	12.00	129	10016	4.55 ug/L	92
62) 1,2-Dibromoethane	12.29	107	12495	4.69 ug/L	100
64) Chlorobenzene	12.89	112	29860	4.76 ug/L	94
65) 1,1,1,2-Tetrachloroethane	12.96	131	8263	4.47 ug/L	92
66) Ethylbenzene	12.98	91	52834	4.63 ug/L	97
67) m,p-Xylene	13.09	106	36804	9.36 ug/L	100
68) o-Xylene	13.72	106	17287	4.68 ug/L	98
69) Styrene	13.77	104	31282	4.64 ug/L	97
70) Bromoform	14.24	173	6995	4.27 ug/L	96
71) Isopropylbenzene	14.25	105	40109	4.69 ug/L	98
72) 1,1,2,2-Tetrachloroethane	14.49	83	17967	4.78 ug/L	100
74) 1,2,3-Trichloropropane	14.70	75	14194	4.83 ug/L	99
75) trans-1,4-Dichloro-2-buten	14.82	89	3797	4.01 ug/L	89
76) n-Propylbenzene	14.83	91	54388	4.65 ug/L	100
77) Bromobenzene	14.88	156	13187	4.72 ug/L	96
78) 1,3,5-Trimethylbenzene	15.06	105	32627	4.70 ug/L	97
79) 2-Chlorotoluene	15.07	91	38111	4.67 ug/L	98

(#) = qualifier out of range (m) = manual integration

Data File : C:\HPCHEM\1\DATA\052606-B\B36410SI.D Vial: 4
Acq On : 26 May 2006 11:22 am Operator:
Sample : 5 PPB Inst : Instrumen
Misc : 5000 Multiplr: 1.00
Integration Params: rteint.p
Quant Time: May 30 8:17 2006 Quant Results File: V805266B.RES

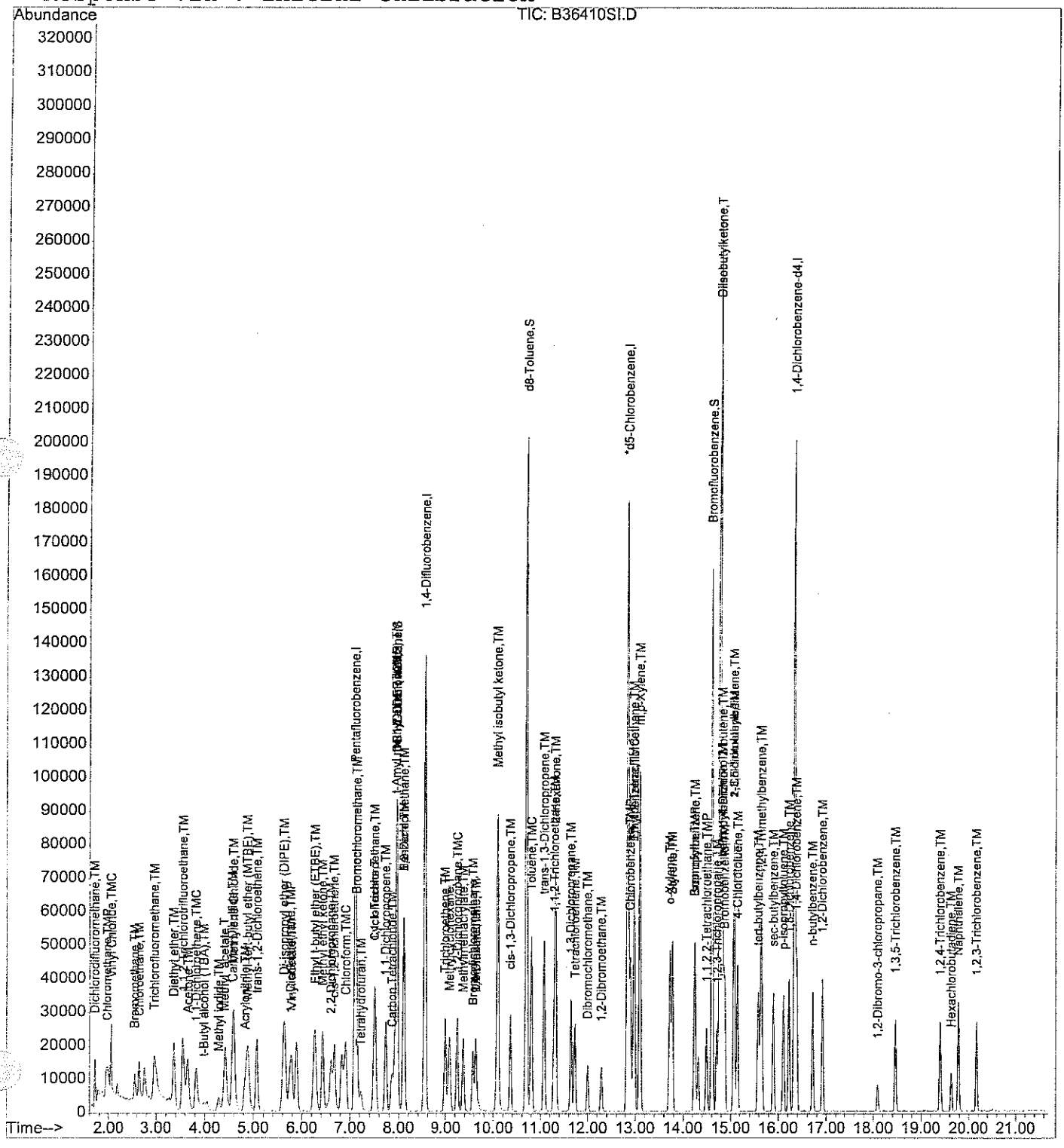
Quant Method : C:\HPCHEM\1\METHODS\V805266B.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Tue May 30 08:15:26 2006
Response via : Initial Calibration
DataAcq Meth : V805266B

	Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
80)	4-Chlorotoluene	15.15	91	35408	4.74	ug/L	97
81)	tert-butylbenzene	15.59	119	23984	4.69	ug/L	100
82)	1,2,4-Trimethylbenzene	15.64	105	32940	4.52	ug/L	97
83)	Diisobutylketone	14.77	57	184422	22.81	ug/L	97
84)	sec-butylbenzene	15.90	105	35578	4.73	ug/L	97
85)	p-isopropyltoluene	16.11	119	27446	4.75	ug/L	98
86)	1,3-Dichlorobenzene	16.23	146	21913	4.73	ug/L	95
88)	1,4-Dichlorobenzene	16.39	146	23363	4.77	ug/L	99
89)	n-butylbenzene	16.71	91	29066	4.75	ug/L	99
90)	1,2-Dichlorobenzene	16.93	146	21570	4.74	ug/L	96
92)	1,2-Dibromo-3-chloropropan	18.09	75	2747	4.68	ug/L	98
94)	1,2,4-Trichlorobenzene	19.40	180	10658	4.68	ug/L	95
95)	Hexachlorobutadiene	19.63	225	2131	4.99	ug/L	92
96)	Naphthalene	19.78	128	33036	4.72	ug/L	100
97)	1,2,3-Trichlorobenzene	20.16	180	10350	4.73	ug/L	95
98)	1,3,5-Trichlorobenzene	18.46	180	11167	4.67	ug/L	97

(#) = qualifier out of range (m) = manual integration

Data File : C:\HPCHEM\1\DATA\052606-B\B36410SI.D Vial: 4
Acq On : 26 May 2006 11:22 am Operator:
Sample : 5 PPB Inst : Instrumen
Misc : 5000 Multiplr: 1.00
MS Integration Params: rteint.p
Quant Time: May 30 8:17 2006 Quant Results File: V805266B.RES

Method : C:\HPCHEM\1\METHODS\V805266B.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Tue May 30 08:15:26 2006
Response via : Initial Calibration



Quantitation Report (Not Reviewed)

Data File : C:\HPCHEM\1\DATA\052606-B\B36411SI.D Vial: 5
Acq On : 26 May 2006 11:52 am Operator:
Sample : 20 PPB Inst : Instrumen
Misc : 5000 Multiplr: 1.00
Integration Params: rteint.p
Quant Time: May 30 8:17 2006 Quant Results File: V805266B.RES

Quant Method : C:\HPCHEM\1\METHODS\V805266B.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Tue May 30 08:15:26 2006
Response via : Initial Calibration
DataAcq Meth : V805266B

5-30-06 SL ✓ 5-30-06

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.11	168	99516	20.00	ug/L	0.00
39) 1,4-Difluorobenzene	8.54	114	154615	20.00	ug/L	0.00
63) *d5-Chlorobenzene	12.84	117	134329	20.00	ug/L	0.00
87) 1,4-Dichlorobenzene-d4	16.34	152	75097	20.00	ug/L	0.00
System Monitoring Compounds						
28) DB HOLDER #28	7.95	65	80205	20.11	ug/L	0.00
Spiked Amount 20.000	Range 70 - 130		Recovery = 100.55%			
36) d4-1,2-Dichloroethane	7.95	65	80205	20.11	ug/L	0.00
Spiked Amount 20.000	Range 70 - 130		Recovery = 100.55%			
54) d8-Toluene	10.69	98	198841	20.08	ug/L	0.00
Spiked Amount 20.000	Range 70 - 130		Recovery = 100.40%			
73) Bromofluorobenzene	14.60	95	74656	20.14	ug/L	0.00
Spiked Amount 20.000	Range 70 - 130		Recovery = 100.70%			
Target Compounds						Qvalue
2) Dichlorodifluoromethane	1.70	85	68427	20.15	ug/L	100
3) Chloromethane	1.97	50	112353	18.62	ug/L	100
4) Vinyl Chloride	2.05	62	58250	17.61	ug/L	100
5) Bromomethane	2.54	94	44450	21.05	ug/L	100
6) Chloroethane	2.63	64	60051	19.40	ug/L	100
7) t-Butyl alcohol (TBA)	3.99	59	32446	104.73	ug/L	100
8) Trichlorofluoromethane	2.96	101	117443	20.03	ug/L	100
9) Diethyl ether	3.36	74	51955	22.34	ug/L	100
10) 1,1,2-Trichlorotrifluoroet	3.56	101	51350	20.21	ug/L	100
12) Acetone	3.64	43	152837	101.34	ug/L	100
13) 1,1-Dichloroethene	3.81	96	47729	20.25	ug/L	100
14) Methyl iodide	4.28	142	56897	18.50	ug/L	100
15) Di-isopropyl ether (DIPE)	5.63	45	245453	19.86	ug/L	100
16) Methylene Chloride	4.59	84	62680	19.08	ug/L	100
17) Carbon Disulfide	4.56	76	185956	19.93	ug/L	100
18) Acrylonitrile	4.81	53	33585	21.24	ug/L	100
19) Methyl-tert-butyl ether (M	4.87	73	177713	20.35	ug/L	100
20) trans-1,2-Dichloroethene	5.06	96	54491	19.79	ug/L	100
21) 1,1-Dichloroethane	5.75	63	108491	19.67	ug/L	100
22) Vinyl acetate	5.78	86	10331	18.91	ug/L	100
23) Methyl ethyl ketone	6.42	43	241712	107.32	ug/L	100
24) Ethyl t-butyl ether (ETBE)	6.26	59	203208	20.25	ug/L	100
25) 2,2-Dichloropropane	6.60	77	71561	20.44	ug/L	100
26) cis-1,2-Dichloroethene	6.67	96	58348	20.05	ug/L	100
27) t-Amyl methyl ether (TAME)	7.93	73	171903	20.39	ug/L	100
29) Chloroform	6.91	83	106006	19.75	ug/L	100

(#) = qualifier out of range (m) = manual integration

Quantitation Report (Not Reviewed)

Data File : C:\HPCHEM\1\DATA\052606-B\B36411SI.D Vial: 5
Acq On : 26 May 2006 11:52 am Operator:
Sample : 20 PPB Inst : Instrumen
Misc : 5000 Multiplr: 1.00
Integration Params: rteint.p
Quant Time: May 30 8:17 2006 Quant Results File: V805266B.RES

Quant Method : C:\HPCHEM\1\METHODS\V805266B.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Tue May 30 08:15:26 2006
Response via : Initial Calibration
DataAcq Meth : V805266B

	Compound	R.T.	QIon	Response	Conc	Unit		Qvalue
30)	Bromochloromethane	7.15	128	26160	20.11	ug/L		100
31)	Tetrahydrofuran	7.21	71	9146	20.64	ug/L		100
32)	1,1,1-Trichloroethane	7.49	97	75805	19.96	ug/L		100
34)	1,1-Dichloropropene	7.73	75	81336	19.92	ug/L		100
35)	Carbon Tetrachloride	7.86	119	50440	20.59	ug/L		100
37)	1,2-Dichloroethane	8.09	62	97386	19.34	ug/L		100
38)	Benzene	8.10	78	227108	19.71	ug/L		100
40)	Trichloroethene	8.99	95	55290	19.61	ug/L		100
41)	1,2-Dichloropropane	9.25	63	61801	20.16	ug/L		100
42)	Methylmethacrylate	9.36	69	51185	22.09	ug/L		100
43)	Methyl acetate	4.38	43	73249	21.23	ug/L		100
44)	Cyclohexane	7.50	56	85258	20.35	ug/L		100
45)	Bromodichloromethane	9.57	83	73826	20.33	ug/L		100
46)	Dibromomethane	9.63	93	45977	20.56	ug/L		100
47)	1,4-Dioxane	9.62	88	21941	556.20	ug/L		100
48)	Methylcyclohexane	9.08	83	60319	20.56	ug/L		100
50)	2-Chloroethylvinyl ether	10.08	63	528	18.19	ug/L	#	1
51)	2-Hexanone	11.32	58	163954	110.16	ug/L		100
52)	Methyl isobutyl ketone	10.08	58	163418	106.74	ug/L		100
53)	cis-1,3-Dichloropropene	10.35	75	92666	20.35	ug/L		100
55)	Toluene	10.80	92	129589	19.68	ug/L		100
56)	trans-1,3-Dichloropropene	11.07	75	85571	20.55	ug/L		100
58)	1,1,2-Trichloroethane	11.29	83	48406	20.13	ug/L		100
59)	1,3-Dichloropropane	11.64	76	96948	20.26	ug/L		100
60)	Tetrachloroethene	11.73	166	46696	19.33	ug/L		100
61)	Dibromochloromethane	12.00	129	47100	20.58	ug/L		100
62)	1,2-Dibromoethane	12.28	107	57589	20.78	ug/L		100
64)	Chlorobenzene	12.89	112	128675	19.36	ug/L		100
65)	1,1,1,2-Tetrachloroethane	12.96	131	39805	20.32	ug/L		100
66)	Ethylbenzene	12.98	91	236649	19.55	ug/L		100
67)	m,p-Xylene	13.09	106	159545	38.27	ug/L		100
68)	o-Xylene	13.72	106	77341	19.76	ug/L		100
69)	Styrene	13.77	104	144606	20.23	ug/L		100
70)	Bromoform	14.24	173	36523	21.04	ug/L		100
71)	Isopropylbenzene	14.25	105	179774	19.84	ug/L		100
72)	1,1,2,2-Tetrachloroethane	14.49	83	83498	20.96	ug/L		100
74)	1,2,3-Trichloropropane	14.70	75	63592	20.41	ug/L		100
75)	trans-1,4-Dichloro-2-buten	14.82	89	20435	20.37	ug/L		100
76)	n-Propylbenzene	14.83	91	242180	19.53	ug/L		100
77)	Bromobenzene	14.88	156	59153	19.95	ug/L		100
78)	1,3,5-Trimethylbenzene	15.06	105	145386	19.77	ug/L		100

(#) = qualifier out of range (m) = manual integration

Quantitation Report (Not Reviewed)

Data File : C:\HPCHEM\1\DATA\052606-B\B36411SI.D Vial: 5
Acq On : 26 May 2006 11:52 am Operator:
Sample : 20 PPB Inst : Instrumen
Misc : 5000 Multiplr: 1.00
IS Integration Params: rteint.p
Quant Time: May 30 8:17 2006 Quant Results File: V805266B.RES

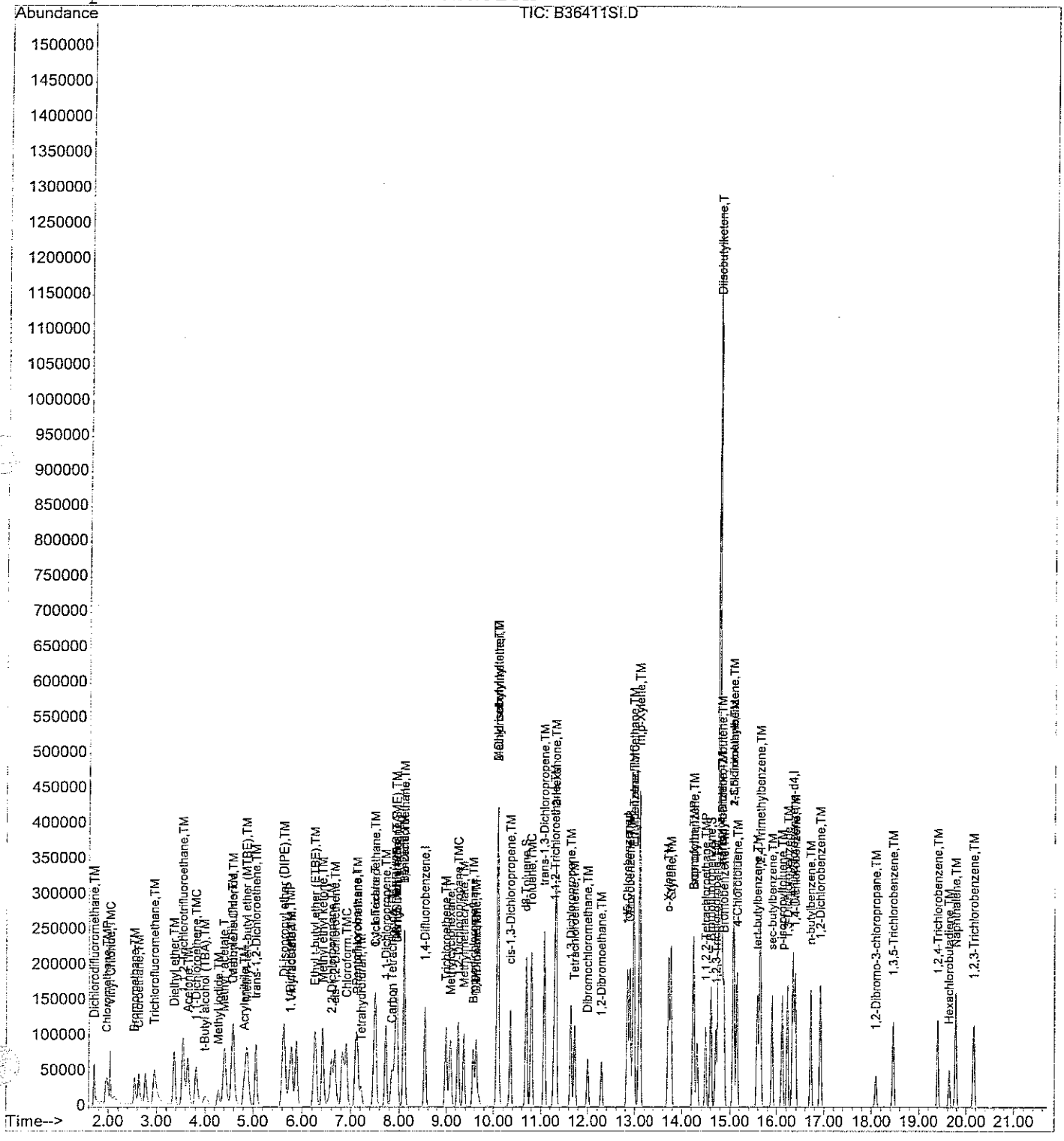
Quant Method : C:\HPCHEM\1\METHODS\V805266B.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Tue May 30 08:15:26 2006
Response via : Initial Calibration
DataAcq Meth : V805266B

	Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
79)	2-Chlorotoluene	15.07	91	170249	19.67	ug/L	100
80)	4-Chlorotoluene	15.15	91	156470	19.76	ug/L	100
81)	tert-butylbenzene	15.59	119	107393	19.79	ug/L	100
82)	1,2,4-Trimethylbenzene	15.64	105	150810	19.53	ug/L	100
83)	Diisobutylketone	14.77	57	872565	101.82	ug/L	100
84)	sec-butylbenzene	15.90	105	158651	19.89	ug/L	100
85)	p-isopropyltoluene	16.11	119	121515	19.85	ug/L	100
86)	1,3-Dichlorobenzene	16.23	146	97419	19.84	ug/L	100
88)	1,4-Dichlorobenzene	16.39	146	102475	19.58	ug/L	100
89)	n-butylbenzene	16.71	91	130203	19.92	ug/L	100
90)	1,2-Dichlorobenzene	16.92	146	95043	19.57	ug/L	100
92)	1,2-Dibromo-3-chloropropan	18.09	75	14162	22.57	ug/L	100
94)	1,2,4-Trichlorobenzene	19.41	180	48769	20.04	ug/L	100
95)	Hexachlorobutadiene	19.63	225	9325	20.45	ug/L	100
96)	Naphthalene	19.78	128	159430	21.32	ug/L	100
97)	1,2,3-Trichlorobenzene	20.16	180	47100	20.17	ug/L	100
98)	1,3,5-Trichlorobenzene	18.46	180	50122	19.62	ug/L	100

(#) = qualifier out of range (m) = manual integration

Data File : C:\HPCHEM\1\DATA\052606-B\B36411SI.D Vial: 5
Acq On : 26 May 2006 11:52 am Operator:
Sample : 20 PPB Inst : Instrumen
Misc : 5000 Multiplr: 1.00
MS Integration Params: rteint.p
Quant Time: May 30 8:17 2006 Quant Results File: V805266B.RES

Method : C:\HPCHEM\1\METHODS\V805266B.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Tue May 30 08:15:26 2006
Response via : Initial Calibration



Quantitation Report (Not Reviewed)

Data File : C:\HPCHEM\1\DATA\052606-B\B36412SI.D Vial: 6
Acq On : 26 May 2006 12:21 pm Operator:
Sample : 50 PPB Inst : Instrumen
Misc : 5000 Multiplr: 1.00
Integration Params: rteint.p
Quant Time: May 30 8:18 2006 Quant Results File: V805266B.RES

Quant Method : C:\HPCHEM\1\METHODS\V805266B.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Tue May 30 08:15:26 2006
Response via : Initial Calibration
DataAcq Meth : V805266B

5.30.06 SL

5.30.06

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.10	168	106866	20.00	ug/L	0.00
39) 1,4-Difluorobenzene	8.54	114	162033	20.00	ug/L	0.00
63) *d5-Chlorobenzene	12.84	117	140743	20.00	ug/L	0.00
87) 1,4-Dichlorobenzene-d4	16.33	152	79036	20.00	ug/L	0.00
System Monitoring Compounds						
28) DB HOLDER #28	7.95	65	81106	18.94	ug/L	0.00
Spiked Amount 20.000	Range 70 - 130		Recovery = 94.70%			
36) d4-1,2-Dichloroethane	7.95	65	81106	18.94	ug/L	0.00
Spiked Amount 20.000	Range 70 - 130		Recovery = 94.70%			
54) d8-Toluene	10.69	98	210368	20.28	ug/L	0.00
Spiked Amount 20.000	Range 70 - 130		Recovery = 101.40%			
73) Bromofluorobenzene	14.61	95	77983	20.08	ug/L	0.00
Spiked Amount 20.000	Range 70 - 130		Recovery = 100.40%			
Target Compounds						Qvalue
2) Dichlorodifluoromethane	1.70	85	180648	49.53	ug/L	100
3) Chloromethane	1.96	50	310904	47.99	ug/L	100
4) Vinyl Chloride	2.04	62	147684	41.58	ug/L	99
5) Bromomethane	2.52	94	155651	68.66	ug/L	99
6) Chloroethane	2.61	64	154458	46.46	ug/L	100
7) t-Butyl alcohol (TBA)	4.02	59	97107	291.90	ug/L	100
8) Trichlorofluoromethane	2.94	101	312112	49.57	ug/L	97
9) Diethyl ether	3.36	74	130563	52.28	ug/L	99
10) 1,1,2-Trichlorotrifluoroet	3.54	101	137967	50.57	ug/L	99
11) Acrolein	3.54	56	210748	107.19	ug/L	98
12) Acetone	3.65	43	406120	250.76	ug/L	100
13) 1,1-Dichloroethene	3.80	96	130376	51.52	ug/L	95
14) Methyl iodide	4.27	142	181815	50.95	ug/L	98
15) Di-isopropyl ether (DIPE)	5.62	45	669121	50.40	ug/L	100
16) Methylene Chloride	4.58	84	167986	47.61	ug/L	99
17) Carbon Disulfide	4.55	76	518054	51.71	ug/L	100
18) Acrylonitrile	4.81	53	90053	53.05	ug/L	100
19) Methyl-tert-butyl ether (M	4.87	73	474963	50.64	ug/L	99
20) trans-1,2-Dichloroethene	5.06	96	146672	49.59	ug/L	97
21) 1,1-Dichloroethane	5.74	63	297617	50.26	ug/L	98
22) Vinyl acetate	5.78	86	31317	51.16	ug/L	99
23) Methyl ethyl ketone	6.42	43	633173	261.78	ug/L	100
24) Ethyl t-butyl ether (ETBE)	6.25	59	543085	50.40	ug/L	100
25) 2,2-Dichloropropane	6.60	77	204253	54.34	ug/L	99
26) cis-1,2-Dichloroethene	6.67	96	159838	51.16	ug/L	99
27) t-Amyl methyl ether (TAME)	7.93	73	463042	51.15	ug/L	99

(#) = qualifier out of range (m) = manual integration

Quantitation Report (Not Reviewed)

Data File : C:\HPCHEM\1\DATA\052606-B\B36412SI.D Vial: 6
Acq On : 26 May 2006 12:21 pm Operator:
Sample : 50 PPB Inst : Instrumen
Misc : 5000 Multiplr: 1.00
Integration Params: rteint.p
Quant Time: May 30 8:18 2006 Quant Results File: V805266B.RES

Quant Method : C:\HPCHEM\1\METHODS\V805266B.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Tue May 30 08:15:26 2006
Response via : Initial Calibration
DataAcq Meth : V805266B

	Compound	R.T.	QIon	Response	Conc	Unit		Qvalue
29)	Chloroform	6.90	83	284680	49.40	ug/L		99
30)	Bromochloromethane	7.14	128	70253	50.29	ug/L		96
31)	Tetrahydrofuran	7.21	71	24451	51.38	ug/L		97
32)	1,1,1-Trichloroethane	7.49	97	205511	50.38	ug/L		97
34)	1,1-Dichloropropene	7.72	75	215356	49.12	ug/L		97
35)	Carbon Tetrachloride	7.85	119	140378	53.37	ug/L		95
37)	1,2-Dichloroethane	8.09	62	257966	47.72	ug/L		99
38)	Benzene	8.10	78	607130	49.07	ug/L		99
40)	Trichloroethene	8.99	95	150214	50.83	ug/L		99
41)	1,2-Dichloropropane	9.24	63	163873	51.02	ug/L		100
42)	Methylmethacrylate	9.36	69	141076	58.10	ug/L		99
43)	Methyl acetate	4.38	43	198594	54.93	ug/L		98
44)	Cyclohexane	7.50	56	226515	51.58	ug/L		100
45)	Bromodichloromethane	9.57	83	200020	52.55	ug/L		98
46)	Dibromomethane	9.64	93	123762	52.80	ug/L		96
47)	1,4-Dioxane	9.62	88	61536	1488.52	ug/L		97
48)	Methylcyclohexane	9.07	83	162944	53.00	ug/L		98
50)	2-Chloroethylvinyl ether	10.09	63	1550	50.97	ug/L	#	1
51)	2-Hexanone	11.32	58	450942	289.10	ug/L		98
52)	Methyl isobutyl ketone	10.08	58	453113	282.42	ug/L		100
53)	cis-1,3-Dichloropropene	10.36	75	258208	54.12	ug/L		100
55)	Toluene	10.80	92	355011	51.44	ug/L		99
56)	trans-1,3-Dichloropropene	11.07	75	247750	56.78	ug/L		100
58)	1,1,2-Trichloroethane	11.28	83	133292	52.89	ug/L		96
59)	1,3-Dichloropropane	11.65	76	262790	52.40	ug/L		99
60)	Tetrachloroethene	11.73	166	130973	51.73	ug/L		97
61)	Dibromochloromethane	12.00	129	136637	56.97	ug/L		100
62)	1,2-Dibromoethane	12.29	107	156352	53.82	ug/L		98
64)	Chlorobenzene	12.89	112	357481	51.35	ug/L		99
65)	1,1,1,2-Tetrachloroethane	12.96	131	115540	56.29	ug/L		98
66)	Ethylbenzene	12.98	91	657066	51.82	ug/L		100
67)	m,p-Xylene	13.10	106	459458	105.18	ug/L		96
68)	o-Xylene	13.72	106	214818	52.38	ug/L		96
69)	Styrene	13.77	104	407077	54.35	ug/L		99
70)	Bromoform	14.24	173	115282	63.38	ug/L		96
71)	Isopropylbenzene	14.24	105	502514	52.94	ug/L		97
72)	1,1,2,2-Tetrachloroethane	14.49	83	227528	54.52	ug/L		100
74)	1,2,3-Trichloropropane	14.70	75	172910	52.98	ug/L		94
75)	trans-1,4-Dichloro-2-buten	14.82	89	62846	59.78	ug/L		92
76)	n-Propylbenzene	14.83	91	687841	52.95	ug/L		98
77)	Bromobenzene	14.87	156	166163	53.49	ug/L		97

(#) = qualifier out of range (m) = manual integration

Data File : C:\HPCHEM\1\DATA\052606-B\B36412SI.D Vial: 6
Acq On : 26 May 2006 12:21 pm Operator:
Sample : 50 PPB Inst : Instrumen
Misc : 5000 Multiplr: 1.00
Integration Params: rteint.p
Quant Time: May 30 8:18 2006 Quant Results File: V805266B.RES

Quant Method : C:\HPCHEM\1\METHODS\V805266B.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Tue May 30 08:15:26 2006
Response via : Initial Calibration
DataAcq Meth : V805266B

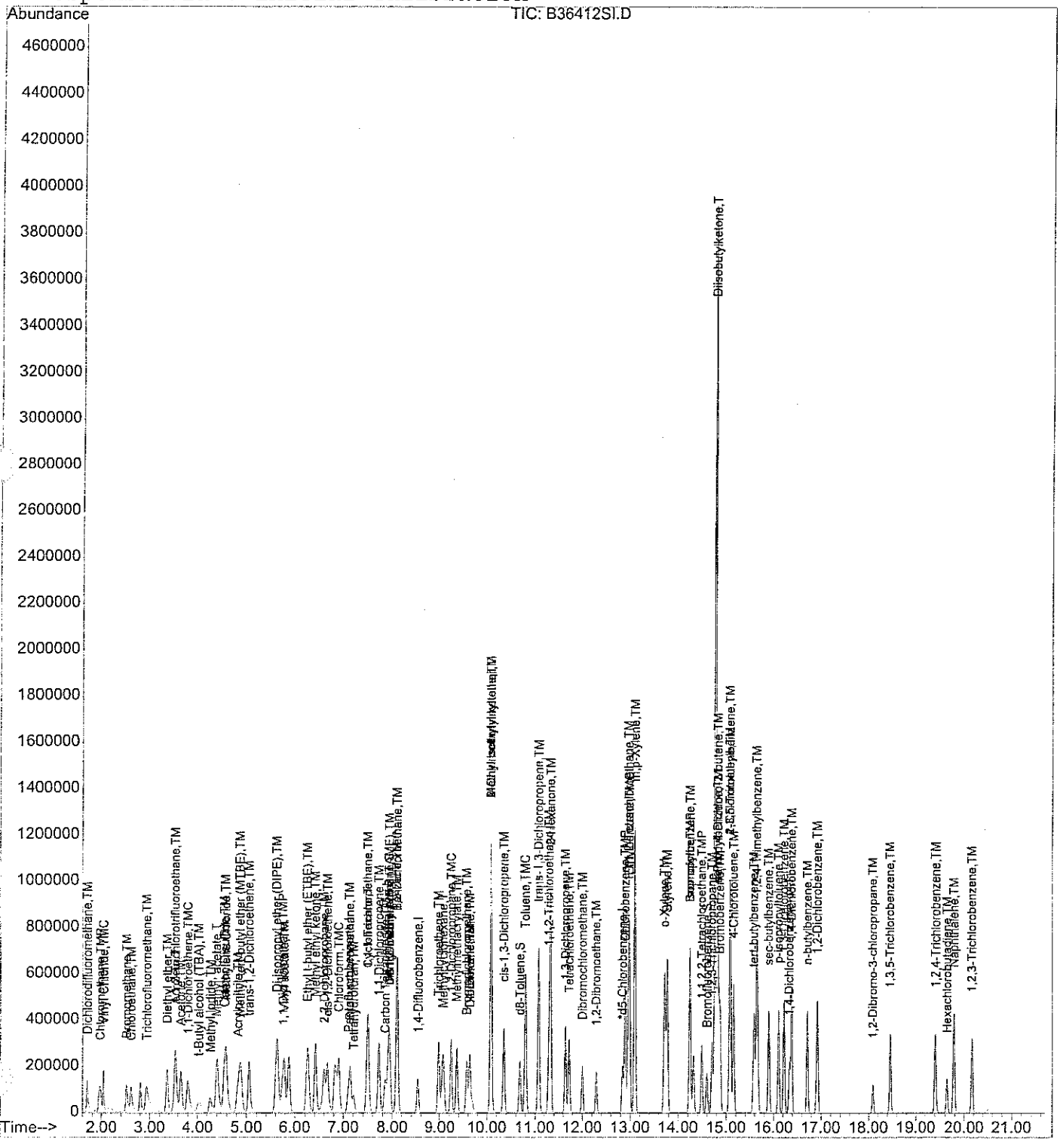
	Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
78)	1,3,5-Trimethylbenzene	15.07	105	409512	53.15	ug/L	96
79)	2-Chlorotoluene	15.08	91	486251	53.63	ug/L	94
80)	4-Chlorotoluene	15.15	91	420225	50.66	ug/L	99
81)	tert-butylbenzene	15.58	119	289674	50.95	ug/L	98
82)	1,2,4-Trimethylbenzene	15.64	105	428352	52.95	ug/L	100
83)	Diisobutylketone	14.77	57	2583765	287.77	ug/L	95
84)	sec-butylbenzene	15.90	105	443938	53.11	ug/L	98
85)	p-isopropyltoluene	16.10	119	343554	53.55	ug/L	98
86)	1,3-Dichlorobenzene	16.22	146	272476	52.97	ug/L	99
88)	1,4-Dichlorobenzene	16.38	146	284569	51.66	ug/L	98
89)	n-butylbenzene	16.72	91	367987	53.48	ug/L	100
90)	1,2-Dichlorobenzene	16.92	146	269534	52.73	ug/L	97
92)	1,2-Dibromo-3-chloropropan	18.09	75	39844	60.32	ug/L	99
94)	1,2,4-Trichlorobenzene	19.41	180	139419	54.42	ug/L	95
95)	Hexachlorobutadiene	19.64	225	26588	55.39	ug/L	98
96)	Naphthalene	19.78	128	445570	56.62	ug/L	100
97)	1,2,3-Trichlorobenzene	20.16	180	135349	55.07	ug/L	99
98)	1,3,5-Trichlorobenzene	18.45	180	142011	52.82	ug/L	99

(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : C:\HPCHEM\1\DATA\052606-B\B36412SI.D Vial: 6
Acq On : 26 May 2006 12:21 pm Operator:
Sample : 50 PPB Inst : Instrumen
Misc : 5000 Multiplr: 1.00
MS Integration Params: rteint.p
Quant Time: May 30 8:18 2006 Quant Results File: V805266B.RES

Method : C:\HPCHEM\1\METHODS\V805266B.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Tue May 30 08:15:26 2006
Response via : Initial Calibration



Quantitation Report (Not Reviewed)

Data File : C:\HPCHEM\1\DATA\052606-B\B36413SI.D Vial: 7
Acq On : 26 May 2006 12:50 pm Operator:
Sample : 100 PPB Inst : Instrumen
Misc : 5000 Multiplr: 1.00
Integration Params: rteint.p
Quant Time: May 30 8:18 2006 Quant Results File: V805266B.RES

Quant Method : C:\HPCHEM\1\METHODS\V805266B.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Tue May 30 08:15:26 2006
Response via : Initial Calibration
DataAcq Meth : V805266B

5-30-06 5-30-06

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.10	168	111566	20.00	ug/L	-0.01
39) 1,4-Difluorobenzene	8.54	114	167116	20.00	ug/L	0.00
63) *d5-Chlorobenzene	12.84	117	147539	20.00	ug/L	0.00
87) 1,4-Dichlorobenzene-d4	16.34	152	84904	20.00	ug/L	0.00
System Monitoring Compounds						
28) DB HOLDER #28	7.95	65	85673	19.16	ug/L	0.00
Spiked Amount 20.000	Range 70 - 130		Recovery = 95.80%			
36) d4-1,2-Dichloroethane	7.95	65	85673	19.16	ug/L	0.00
Spiked Amount 20.000	Range 70 - 130		Recovery = 95.80%			
54) d8-Toluene	10.68	98	214597	20.05	ug/L	0.00
Spiked Amount 20.000	Range 70 - 130		Recovery = 100.25%			
73) Bromofluorobenzene	14.60	95	80470	19.77	ug/L	0.00
Spiked Amount 20.000	Range 70 - 130		Recovery = 98.85%			
Target Compounds						Qvalue
2) Dichlorodifluoromethane	1.69	85	365750	96.06	ug/L	100
3) Chloromethane	1.97	50	612175	90.52	ug/L	100
4) Vinyl Chloride	2.04	62	285724	77.06	ug/L	99
5) Bromomethane	2.49	94	329831	139.36	ug/L	100
6) Chloroethane	2.59	64	301072	86.75	ug/L	100
7) t-Butyl alcohol (TBA)	4.07	59	189319	545.11	ug/L	99
8) Trichlorofluoromethane	2.91	101	627815	95.51	ug/L	99
9) Diethyl ether	3.37	74	244035	93.59	ug/L	97
10) 1,1,2-Trichlorotrifluoroet	3.53	101	284585	99.91	ug/L	98
11) Acrolein	3.54	56	418402	203.84	ug/L	98
12) Acetone	3.67	43	784698	464.10	ug/L	100
13) 1,1-Dichloroethene	3.78	96	275220	104.17	ug/L	98
14) Methyl iodide	4.26	142	388929	99.83	ug/L	98
15) Di-isopropyl ether (DIPE)	5.63	45	1399671	100.99	ug/L	98
16) Methylene Chloride	4.57	84	346374	94.03	ug/L	98
17) Carbon Disulfide	4.53	76	1093552	104.57	ug/L	100
18) Acrylonitrile	4.82	53	181043	102.15	ug/L	99
19) Methyl-tert-butyl ether (M	4.88	73	978488	99.93	ug/L	98
20) trans-1,2-Dichloroethene	5.05	96	309681	100.30	ug/L	96
21) 1,1-Dichloroethane	5.73	63	627237	101.46	ug/L	98
22) Vinyl acetate	5.78	86	64436	99.65	ug/L	96
23) Methyl ethyl ketone	6.42	43	1235494	489.29	ug/L	99
24) Ethyl t-butyl ether (ETBE)	6.26	59	1147004	101.97	ug/L	99
25) 2,2-Dichloropropane	6.59	77	427783	109.02	ug/L	99
26) cis-1,2-Dichloroethene	6.66	96	334178	102.45	ug/L	99
27) t-Amyl methyl ether (TAME)	7.93	73	961715	101.76	ug/L	98

(#) = qualifier out of range (m) = manual integration

Quantitation Report (Not Reviewed)

Data File : C:\HPCHEM\1\DATA\052606-B\B36413SI.D Vial: 7
Acq On : 26 May 2006 12:50 pm Operator:
Sample : 100 PPB Inst : Instrumen
Misc : 5000 Multiplr: 1.00
IS Integration Params: rteint.p
Quant Time: May 30 8:18 2006 Quant Results File: V805266B.RES

Quant Method : C:\HPCHEM\1\METHODS\V805266B.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Tue May 30 08:15:26 2006
Response via : Initial Calibration
DataAcq Meth : V805266B

	Compound	R.T.	QIon	Response	Conc	Unit		Qvalue
29)	Chloroform	6.91	83	595196	98.93	ug/L		99
30)	Bromochloromethane	7.13	128	145869	100.01	ug/L		97
31)	Tetrahydrofuran	7.21	71	48112	96.83	ug/L		93
32)	1,1,1-Trichloroethane	7.49	97	438220	102.91	ug/L		97
34)	1,1-Dichloropropene	7.72	75	451450	98.63	ug/L		97
35)	Carbon Tetrachloride	7.84	119	304079	110.73	ug/L		94
37)	1,2-Dichloroethane	8.09	62	529621	93.84	ug/L		100
38)	Benzene	8.10	78	1295091	100.26	ug/L		100
40)	Trichloroethene	8.99	95	313844	102.96	ug/L		99
41)	1,2-Dichloropropane	9.25	63	341999	103.23	ug/L		100
42)	Methylmethacrylate	9.36	69	283494	113.20	ug/L		97
43)	Methyl acetate	4.39	43	398520	106.87	ug/L		97
44)	Cyclohexane	7.49	56	480579	106.10	ug/L		98
45)	Bromodichloromethane	9.57	83	422995	107.76	ug/L		98
46)	Dibromomethane	9.63	93	249793	103.32	ug/L		97
47)	1,4-Dioxane	9.63	88	120958	2836.90	ug/L		97
48)	Methylcyclohexane	9.07	83	339713	107.15	ug/L		97
50)	2-Chloroethylvinyl ether	10.08	63	3359	107.09	ug/L	#	1
51)	2-Hexanone	11.33	58	894879	556.27	ug/L		94
52)	Methyl isobutyl ketone	10.08	58	912974	551.74	ug/L		99
53)	cis-1,3-Dichloropropene	10.35	75	550104	111.79	ug/L		100
55)	Toluene	10.80	92	754229	105.97	ug/L		100
56)	trans-1,3-Dichloropropene	11.07	75	537353	119.41	ug/L		98
58)	1,1,2-Trichloroethane	11.29	83	275080	105.83	ug/L		96
59)	1,3-Dichloropropane	11.64	76	546361	105.63	ug/L		98
60)	Tetrachloroethene	11.73	166	277665	106.33	ug/L		97
61)	Dibromochloromethane	12.01	129	293471	118.65	ug/L		100
62)	1,2-Dibromoethane	12.29	107	323866	108.09	ug/L		98
64)	Chlorobenzene	12.89	112	770807	105.61	ug/L		98
65)	1,1,1,2-Tetrachloroethane	12.96	131	257300	119.58	ug/L		97
66)	Ethylbenzene	12.98	91	1426808	107.34	ug/L		98
67)	m,p-Xylene	13.09	106	1035626	226.15	ug/L		94
68)	o-Xylene	13.72	106	470031	109.34	ug/L		99
69)	Styrene	13.77	104	879012	111.96	ug/L		99
70)	Bromoform	14.23	173	255951	134.24	ug/L		96
71)	Isopropylbenzene	14.24	105	1110727	111.62	ug/L		97
72)	1,1,2,2-Tetrachloroethane	14.49	83	452558	103.44	ug/L		100
74)	1,2,3-Trichloropropane	14.71	75	348327	101.81	ug/L		98
75)	trans-1,4-Dichloro-2-buten	14.82	89	139915	126.95	ug/L	#	74
76)	n-Propylbenzene	14.84	91	1524553	111.96	ug/L		98
77)	Bromobenzene	14.88	156	358162	109.99	ug/L		98

(#) = qualifier out of range (m) = manual integration

Quantitation Report (Not Reviewed)

Data File : C:\HPCHEM\1\DATA\052606-B\B36413SI.D Vial: 7
Acq On : 26 May 2006 12:50 pm Operator:
Sample : 100 PPB Inst : Instrumen
Misc : 5000 Multiplr: 1.00
RTE Integration Params: rteint.p
Quant Time: May 30 8:18 2006 Quant Results File: V805266B.RES

Quant Method : C:\HPCHEM\1\METHODS\V805266B.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Tue May 30 08:15:26 2006
Response via : Initial Calibration
DataAcq Meth : V805266B

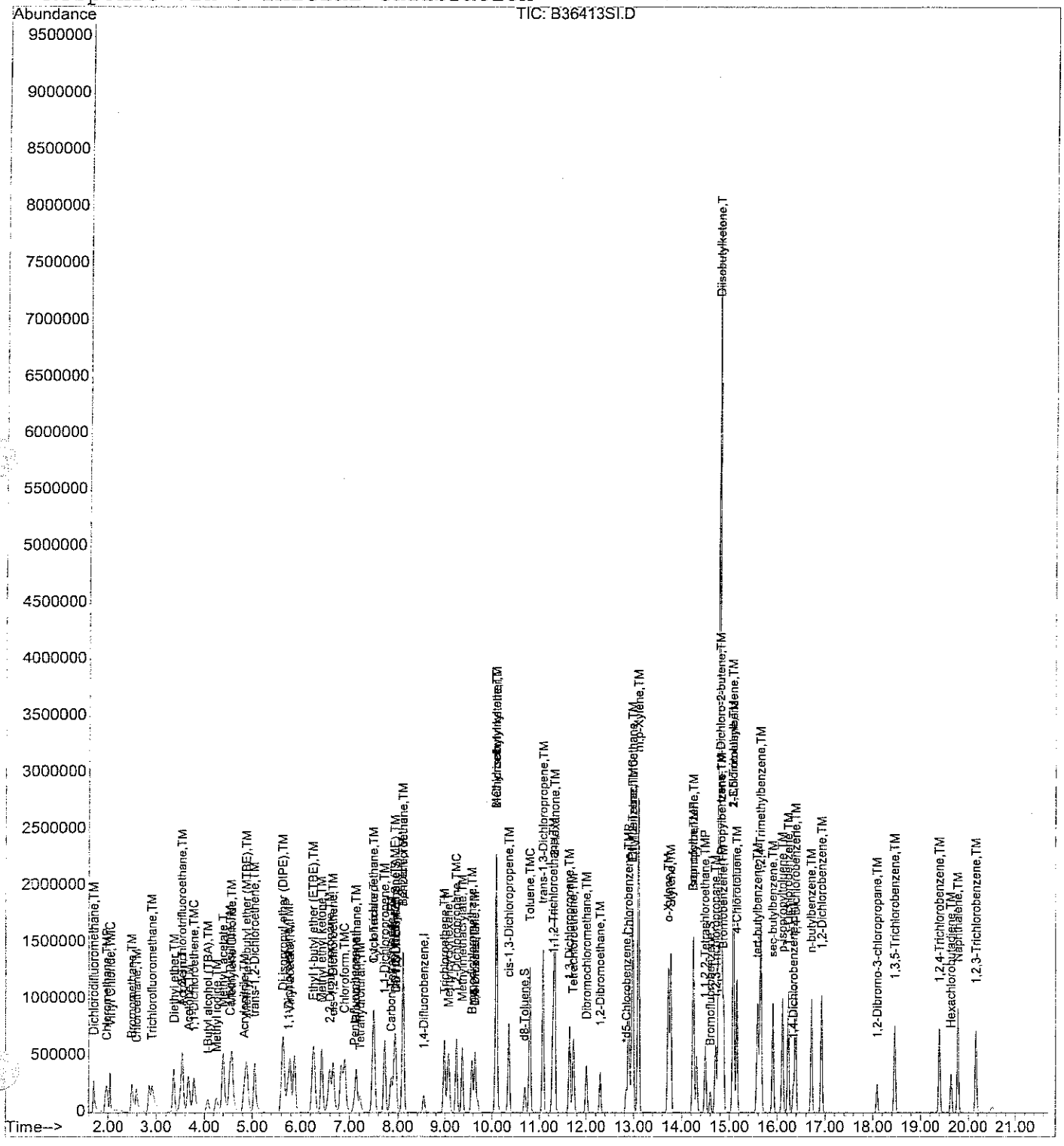
	Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
78)	1,3,5-Trimethylbenzene	15.07	105	919894	113.89	ug/L	95
79)	2-Chlorotoluene	15.07	91	1049433	110.41	ug/L	97
80)	4-Chlorotoluene	15.15	91	942525	108.38	ug/L	99
81)	tert-butylbenzene	15.59	119	645412	108.30	ug/L	98
82)	1,2,4-Trimethylbenzene	15.64	105	954078	112.51	ug/L	99
83)	Diisobutylketone	14.78	57	5488798	583.15	ug/L	87
84)	sec-butylbenzene	15.90	105	964250	110.05	ug/L	100
85)	p-isopropyltoluene	16.11	119	758633	112.80	ug/L	99
86)	1,3-Dichlorobenzene	16.23	146	594822	110.31	ug/L	99
88)	1,4-Dichlorobenzene	16.38	146	626234	105.83	ug/L	98
89)	n-butylbenzene	16.71	91	804893	108.89	ug/L	99
90)	1,2-Dichlorobenzene	16.93	146	590893	107.61	ug/L	97
92)	1,2-Dibromo-3-chloropropan	18.09	75	80755	113.81	ug/L	94
94)	1,2,4-Trichlorobenzene	19.41	180	314525	114.29	ug/L	95
95)	Hexachlorobutadiene	19.64	225	61069	118.43	ug/L	99
96)	Naphthalene	19.78	128	949958	112.36	ug/L	100
97)	1,2,3-Trichlorobenzene	20.16	180	299349	113.38	ug/L	98
98)	1,3,5-Trichlorobenzene	18.46	180	318812	110.38	ug/L	100

(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : C:\HPCHEM\1\DATA\052606-B\B36413SI.D Vial: 7
Acq On : 26 May 2006 12:50 pm Operator:
Sample : 100 PPB Inst : Instrumen
Misc : 5000 Multiplr: 1.00
MS Integration Params: rteint.p
Quant Time: May 30 8:18 2006 Quant Results File: V805266B.RES

Method : C:\HPCHEM\1\METHODS\V805266B.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Tue May 30 08:15:26 2006
Response via : Initial Calibration



Evaluate Continuing Calibration Report

Data File : C:\HPCHEM\1\DATA\053006-B\B36459SC.D Vial: 3
Acq On : 30 May 2006 9:35 am Operator:
Sample : 20 PPB Inst : Instrumen
Misc : 5000 Multiplr: 1.00
MS Integration Params: rteint.p

Method : C:\HPCHEM\1\METHODS\V805266B.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Tue May 30 08:22:32 2006
Response via : Multiple Level Calibration

Min. RRF : 0.100 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
Max. RRF Dev : 20% Max. Rel. Area : 200%

		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
1	I	Pentafluorobenzene	1.000	1.000	0.0	123	0.00
2	TM	Dichlorodifluoromethane	0.683	0.692	-1.3	124	0.00
3	TMP	Chloromethane	1.212	1.169	3.5	127	-0.03
4	TMC	Vinyl Chloride	0.665	0.540	18.8	114	0.00
5	TM	Bromomethane	0.424	0.348	17.9	96	0.00
6	TM	Chloroethane	0.622	0.558	10.3	114	0.00
7	TM	t-Butyl alcohol (TBA)	0.062	0.070#	-12.9	132	0.00
8	TM	Trichlorofluoromethane	1.178	1.113	5.5	116	0.00
9	TM	Diethyl ether	0.467	0.465	0.4	110	0.00
10	TM	1,1,2-Trichlorotrifluoroeth	0.511	0.508	0.6	121	0.00
11	T	Acrolein	0.368	0.343	6.8	109	0.00
12	TM	Acetone	0.303	0.278	8.3	111	0.00
13	TMC	1,1-Dichloroethene	0.474	0.472	0.4	121	0.00
14	TM	Methyl iodide	0.462	0.421	8.9	91	0.00
15	TM	Di-isopropyl ether (DIPE)	2.484	2.447	1.5	122	0.00
16	TM	Methylene Chloride	0.660	0.637	3.5	124	0.00
17	TM	Carbon Disulfide	1.875	1.842	1.8	121	0.00
18	TM	Acrylonitrile	0.318	0.316	0.6	115	0.00
19	TM	Methyl-tert-butyl ether (MT	1.755	1.768	-0.7	122	0.00
20	TM	trans-1,2-Dichloroethene	0.553	0.538	2.7	121	0.00
21	TMP	1,1-Dichloroethane	1.108	1.075	3.0	121	0.00
22	TM	Vinyl acetate	0.084	0.114	NT -35.7#	135	0.00
23	TM	Methyl ethyl ketone	0.453	0.431	4.9	109	0.00
24	TM	Ethyl t-butyl ether (ETBE)	2.017	2.032	-0.7	122	0.00
25	TM	2,2-Dichloropropane	0.703	0.801	-13.9	137	0.00
26	TM	cis-1,2-Dichloroethene	0.585	0.577	1.4	121	0.00
27	TM	t-Amyl methyl ether (TAME)	1.694	1.718	-1.4	122	0.00
28	S	DB HOLDER #28	0.802	0.775	3.4	118	0.00
29	TMC	Chloroform	1.079	1.035	4.1	120	0.00
30	TM	Bromochloromethane	0.261	0.257	1.5	121	0.00
31	TM	Tetrahydrofuran	0.089	0.084#	5.6	113	0.00
32	TM	1,1,1-Trichloroethane	0.763	0.788	-3.3	127	0.00
33	T	NO COMPOUND	0.000	0.000#	0.0	0#	-0.13
34	TM	1,1-Dichloropropene	0.821	0.794	3.3	120	0.00
35	TM	Carbon Tetrachloride	0.492	0.549	-11.6	133	0.00
36	S	d4-1,2-Dichloroethane	0.802	0.775	3.4	118	0.00
37	TM	1,2-Dichloroethane	1.012	0.954	5.7	120	0.00
38	TM	Benzene	2.316	2.213	4.4	119	0.00
39	I	1,4-Difluorobenzene	1.000	1.000	0.0	123	0.00
40	TM	Trichloroethene	0.365	0.359	1.6	123	0.00

(#) = Out of Range

Evaluate Continuing Calibration Report

Data File : C:\HPCHEM\1\DATA\053006-B\B36459SC.D Vial: 3
Acq On : 30 May 2006 9:35 am Operator:
Sample : 20 PPB Inst : Instrumen
Misc : 5000 Multiplr: 1.00
MS Integration Params: rteint.p

Method : C:\HPCHEM\1\METHODS\V805266B.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Tue May 30 08:22:32 2006
Response via : Multiple Level Calibration

Min. RRF : 0.100 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
Max. RRF Dev : 20% Max. Rel. Area : 200%

		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
41	TMC	1,2-Dichloropropane	0.396	0.390	1.5	120	0.00
42	TM	Methylmethacrylate	0.300	0.319	-6.3	118	0.00
43	T	Methyl acetate	0.446	0.456	-2.2	118	0.00
44	T	Cyclohexane	0.542	0.541	0.2	121	0.00
45	TM	Bromodichloromethane	0.470	0.477	-1.5	123	0.00
46	TM	Dibromomethane	0.289	0.291	-0.7	120	0.00
47	TM	1,4-Dioxane	0.005	0.005#	0.0	117	0.00
48	T	Methylcyclohexane	0.379	0.383	-1.1	121	0.00
49	T	NO COMPOUND	0.000	0.000#	0.0	0#	-0.13
50	T	2-Chloroethylvinyl ether	0.004	0.003#	25.0#	107	0.00
51	TM	2-Hexanone	0.193	0.190	1.6	110	0.00
52	TM	Methyl isobutyl ketone	0.198	0.195	1.5	113	0.00
53	TM	cis-1,3-Dichloropropene	0.589	0.621	-5.4	127	0.00
54	S	d8-Toluene	1.281	1.266	1.2	121	0.00
55	TMC	Toluene	0.852	0.823	3.4	121	0.00
56	TM	trans-1,3-Dichloropropene	0.539	0.580	-7.6	129	0.00
57	T	NO COMPOUND	0.000	0.000#	0.0	0#	-0.13
58	TM	1,1,2-Trichloroethane	0.311	0.309	0.6	121	0.00
59	TM	1,3-Dichloropropane	0.619	0.608	1.8	119	0.00
60	TM	Tetrachloroethene	0.313	0.312	0.3	127	0.00
61	TM	Dibromochloromethane	0.296	0.319	-7.8	129	0.00
62	TM	1,2-Dibromoethane	0.359	0.365	-1.7	120	0.00
63	I	*d5-Chlorobenzene	1.000	1.000	0.0	121	0.00
64	TMP	Chlorobenzene	0.989	0.976	1.3	124	0.00
65	TM	1,1,1,2-Tetrachloroethane	0.292	0.310	-6.2	127	0.00
66	TMC	Ethylbenzene	1.802	1.738	3.6	120	0.00
67	TM	m,p-Xylene	0.621	0.603	2.9	123	0.00
68	TM	o-Xylene	0.583	0.575	1.4	121	0.00
69	TM	Styrene	1.064	1.066	-0.2	120	0.00
70	TMP	Bromoform	0.258	0.296	-14.7	132	0.00
71	TM	Isopropylbenzene	1.349	1.324	1.9	120	0.00
72	TMP	1,1,2,2-Tetrachloroethane	0.593	0.583	1.7	114	0.00
73	S	Bromofluorobenzene	0.552	0.555	-0.5	121	0.00
74	TM	1,2,3-Trichloropropane	0.464	0.456	1.7	117	-0.02
75	TM	trans-1,4-Dichloro-2-butene	0.149	0.161	-8.1	129	0.00
76	TM	n-Propylbenzene	1.846	1.801	2.4	121	0.00
77	TM	Bromobenzene	0.441	0.445	-0.9	123	0.00
78	TM	1,3,5-Trimethylbenzene	1.095	1.069	2.4	120	0.00
79	TM	2-Chlorotoluene	1.289	1.276	1.0	122	0.00
80	TM	4-Chlorotoluene	1.179	1.159	1.7	121	0.00

(#) = Out of Range

Evaluate Continuing Calibration Report

Data File : C:\HPCHEM\1\DATA\053006-B\B36459SC.D Vial: 3
Acq On : 30 May 2006 9:35 am Operator:
Sample : 20 PPB Inst : Instrumen
Misc : 5000 Multiplr: 1.00
MS Integration Params: rteint.p

Method : C:\HPCHEM\1\METHODS\V805266B.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Tue May 30 08:22:32 2006
Response via : Multiple Level Calibration

Min. RRF : 0.100 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
Max. RRF Dev : 20% Max. Rel. Area : 200%

		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
81	TM	tert-butylbenzene	0.808	0.811	-0.4	123	0.00
82	TM	1,2,4-Trimethylbenzene	1.149	1.121	2.4	121	0.00
83	T	Diisobutylketone	1.276	1.287	-0.9	120	0.00
84	TM	sec-butylbenzene	1.188	1.188	0.0	122	0.00
85	TM	p-isopropyltoluene	0.912	0.924	-1.3	124	0.00
86	TM	1,3-Dichlorobenzene	0.731	0.723	1.1	121	0.00
87	I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	124	0.00
88	TM	1,4-Dichlorobenzene	1.394	1.345	3.5	122	0.00
89	TM	n-butylbenzene	1.741	1.711	1.7	122	0.00
90	TM	1,2-Dichlorobenzene	1.294	1.260	2.6	123	0.00
91	T	NO COMPOUND	0.000	0.000#	0.0	0#	-0.13
92	TM	1,2-Dibromo-3-chloropropane	0.167	0.182	-9.0	119	0.00
93	T	NO COMPOUND	0.000	0.000#	0.0	0#	-0.13
94	TM	1,2,4-Trichlorobenzene	0.648	0.643	0.8	122	0.00
95	TM	Hexachlorobutadiene	0.121	0.134	-10.7	133	0.00
96	TM	Naphthalene	1.991	2.023	-1.6	118	0.00
97	TM	1,2,3-Trichlorobenzene	0.622	0.620	0.3	122	0.00
98	TM	1,3,5-Trichlorobenzene	0.680	0.673	1.0	125	0.00

Evaluate Continuing Calibration Report

Data File : D:\HPCHEM\1\DATA\053006C\C59118SC.D Vial: 3
Acq On : 30 May 2006 8:45 am Operator:
Sample : 20 PPB Inst : Instr_C
Misc : 5000 Multiplr: 1.00
MS Integration Params: rteint.p

Method : D:\HPCHEM\1\METHODS\V805226C.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Mon May 22 14:18:09 2006
Response via : Multiple Level Calibration

5.30.06 SU 6.5.06

Min. RRF : 0.100 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
Max. RRF Dev : 20% Max. Rel. Area : 200%

		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
1	I	Pentafluorobenzene	1.000	1.000	0.0	106	-0.06
2	TM	Dichlorodifluoromethane	0.601	0.617	-2.7	113	-0.02
3	TMP	Chloromethane	0.636	0.582	8.5	101	-0.03
4	TMC	Vinyl Chloride	0.642	0.608	5.3	104	-0.02
5	TM	Bromomethane	0.350	0.346	1.1	92	-0.02
6	TM	Chloroethane	0.347	0.310	10.7	106	-0.02
7	TM	t-Butyl alcohol (TBA)	0.058	0.050#	13.8	99	-0.06
8	TM	Trichlorofluoromethane	1.213	1.175	3.1	108	-0.02
9	TM	Diethyl ether	0.208	0.186	10.6	101	-0.04
10	T	1,1,2-Trichlorotrifluoroeth	0.543	0.488	10.1	101	-0.04
11	T	Acrolein	0.070	0.060#	14.3	84	-0.04
12	TM	Acetone	0.183	0.169	7.7	106	-0.04
13	TMC	1,1-Dichloroethene	0.416	0.355	14.7	98	-0.04
14	T	Methyl iodide	0.731	0.580	20.7#	74	-0.05
15	TM	Di-isopropyl ether (DIPE)	1.824	1.650	9.5	101	-0.06
16	TM	Methylene Chloride	0.431	0.384	10.9	101	-0.06
17	TM	Carbon Disulfide	1.337	1.259	5.8	101	-0.04
18	T	Acrylonitrile	0.188	0.162	13.8	96	-0.05
19	TM	Methyl-tert-butyl ether (MT	1.513	1.369	9.5	103	-0.05
20	TM	trans-1,2-Dichloroethene	0.944	0.856	9.3	105	-0.05
21	TMP	1,1-Dichloroethane	1.081	1.006	6.9	104	-0.05
22	T	Vinyl acetate	1.184	1.150	2.9	110	-0.05
23	TM	Methyl ethyl ketone	0.050	0.044#	12.0	100	-0.06
24	TM	Ethyl t-butyl ether (ETBE)	1.913	1.780	7.0	104	-0.05
25	TM	2,2-Dichloropropane	0.926	1.143	-23.4#	157	-0.04
26	TM	cis-1,2-Dichloroethene	0.612	0.501	18.1	86	-0.05
27	TM	t-Amyl methyl ether (TAME)	1.328	1.229	7.5	103	-0.05
28	S	DB Holder #28	0.782	0.717	8.3	103	-0.05
29	TMC	Chloroform	1.094	1.217	-11.2	131	-0.06
30	TM	Bromochloromethane	0.336	0.311	7.4	101	-0.05
31	TM	Tetrahydrofuran	0.056	0.054#	3.6	104	-0.05
32	TM	1,1,1-Trichloroethane	0.932	0.884	5.2	105	-0.05
33	T	NO COMPOUND	0.000	0.000#	0.0	0#	-0.49
34	TM	1,1-Dichloropropene	0.767	0.727	5.2	105	-0.05
35	TM	Carbon Tetrachloride	0.784	0.782	0.3	106	-0.05
36	S	d4-1,2-Dichloroethane	0.782	0.717	8.3	103	-0.05
37	TM	1,2-Dichloroethane	0.992	0.922	7.1	103	-0.05
38	TM	Benzene	1.924	1.697	11.8	104	-0.05
39	I	1,4-Difluorobenzene	1.000	1.000	0.0	107	-0.04
40	TM	Trichloroethene	0.377	0.346	8.2	102	-0.05

(#) = Out of Range

Evaluate Continuing Calibration Report

Data File : D:\HPCHEM\1\DATA\053006C\C59118SC.D Vial: 3
Acq On : 30 May 2006 8:45 am Operator:
Sample : 20 PPB Inst : Instr_C
Misc : 5000 Multiplr: 1.00
MS Integration Params: rteint.p

Method : D:\HPCHEM\1\METHODS\V805226C.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Mon May 22 14:18:09 2006
Response via : Multiple Level Calibration

5:31 JL

Min. RRF : 0.100 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
Max. RRF Dev : 20% Max. Rel. Area : 200%

		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
41	TMC	1,2-Dichloropropane	0.349	0.325	6.9	104	-0.04
42	T	Methylmethacrylate	0.221	0.200	9.5	100	-0.05
43	T	Methyl acetate	0.283	0.249	12.0	100	-0.05
44	T	Cyclohexane	0.571	0.522	8.6	101	-0.06
45	TM	Bromodichloromethane	0.440	0.424	3.6	106	-0.05
46	TM	Dibromomethane	0.271	0.245	9.6	97	-0.05
47	T	1,4-Dioxane	0.004	0.003#	25.0#	107	-0.05
48	T	Methylcyclohexane	0.483	0.440	8.9	102	-0.05
49	T	NO COMPOUND	0.000	0.000#	0.0	0#	-0.49
50	T	2-Chloroethylvinyl ether	0.107	0.143	33.6#	134	-0.04
51	TM	2-Hexanone	0.128	0.114	10.9	103	-0.05
52	TM	Methyl isobutyl ketone	0.130	0.118	9.2	101	-0.04
53	TM	cis-1,3-Dichloropropene	0.507	0.492	3.0	105	-0.05
54	S	d8-Toluene	1.143	1.015	11.2	103	-0.05
55	TMC	Toluene	0.724	0.662	8.6	101	-0.04
56	TM	trans-1,3-Dichloropropene	0.482	0.457	5.2	106	-0.04
57	T	NO COMPOUND	0.000	0.000#	0.0	0#	-0.49
58	TM	1,1,2-Trichloroethane	0.240	0.223	7.1	103	-0.05
59	TM	1,3-Dichloropropane	0.463	0.432	6.7	104	-0.04
60	TM	Tetrachloroethene	0.435	0.423	2.8	108	-0.04
61	TM	Dibromochloromethane	0.378	0.369	2.4	106	-0.04
62	TM	1,2-Dibromoethane	0.406	0.380	6.4	103	-0.05
63	I	*d5-Chlorobenzene	1.000	1.000	0.0	109	-0.05
64	TMP	Chlorobenzene	1.073	0.989	7.8	104	-0.05
65	TM	1,1,1,2-Tetrachloroethane	0.379	0.358	5.5	106	-0.05
66	TMC	Ethylbenzene	1.554	1.393	10.4	101	-0.05
67	TM	m,p-Xylene	0.623	0.578	7.2	106	-0.04
68	TM	o-Xylene	0.603	0.562	6.8	107	-0.04
69	TM	Styrene	1.061	0.947	10.7	101	-0.05
70	TMP	Bromoform	0.325	0.331	-1.8	113	-0.05
71	TM	Isopropylbenzene	1.606	1.464	8.8	104	-0.05
72	TMP	1,1,2,2-Tetrachloroethane	0.426	0.379	11.0	103	-0.04
73	S	Bromofluorobenzene	0.498	0.442	11.2	105	-0.04
74	TM	1,2,3-Trichloropropane	0.309	0.268	13.3	101	-0.05
75	TM	trans-1,4-Dichloro-2-butene	0.126	0.115	8.7	106	-0.04
76	TM	n-Propylbenzene	1.775	1.573	11.4	101	-0.05
77	TM	Bromobenzene	0.451	0.423	6.2	107	-0.05
78	TM	1,3,5-Trimethylbenzene	1.286	1.162	9.6	103	-0.05
79	TM	2-Chlorotoluene	1.191	1.065	10.6	102	-0.04
80	TM	4-Chlorotoluene	1.097	0.980	10.7	105	-0.05

(#) = Out of Range

Evaluate Continuing Calibration Report

Data File : D:\HPCHEM\1\DATA\053006C\C59118SC.D Vial: 3
Acq On : 30 May 2006 8:45 am Operator:
Sample : 20 PPB Inst : Instr_C
Misc : 5000 Multiplr: 1.00
MS Integration Params: rteint.p

Method : D:\HPCHEM\1\METHODS\V805226C.M (RTE Integrator)
Title : 8260 Purgable Organics
Last Update : Mon May 22 14:18:09 2006
Response via : Multiple Level Calibration

Min. RRF : 0.100 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
Max. RRF Dev : 20% Max. Rel. Area : 200%

		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
81	TM	tert-butylbenzene	1.124	1.044	7.1	107	-0.04
82	TM	1,2,4-Trimethylbenzene	1.319	1.155	12.4	105	-0.04
83	T	Diisobutylketone	0.919	0.784	14.7	97	-0.05
84	TM	sec-butylbenzene	1.770	1.585	10.5	105	-0.04
85	TM	p-isopropyltoluene	1.396	1.281	8.2	106	-0.04
86	TM	1,3-Dichlorobenzene	0.754	0.695	7.8	106	-0.04
87	I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	111	-0.04
88	TM	1,4-Dichlorobenzene	1.592	1.481	7.0	106	-0.04
89	TM	n-butylbenzene	2.667	2.354	11.7	101	-0.04
90	TM	1,2-Dichlorobenzene	1.461	1.365	6.6	104	-0.04
91	T	NO COMPOUND	0.000	0.000#	0.0	0#	-0.49
92	TM	1,2-Dibromo-3-chloropropane	0.149	0.142	4.7	111	-0.04
93	T	NO COMPOUND	0.000	0.000#	0.0	0#	-0.49
94	TM	1,2,4-Trichlorobenzene	1.221	1.217	0.3	114	-0.05
95	TM	Hexachlorobutadiene	0.599	0.611	-2.0	116	-0.04
96	TM	Naphthalene	2.315	2.154	7.0	111	-0.04
97	TM	1,2,3-Trichlorobenzene	1.086	1.077	0.8	115	-0.05
98	T	1,3,5-Trichlorobenzene	1.359	1.335	1.8	113	-0.05

analytics environmental laboratory LLC

VOLATILE ORGAN
INTERNAL STANDARD AREA AND RT SUM

Project #: 218736.002 Quote #: SDG#: 56637

BFB File ID: C59021T.D Site ID: GP Old Town Mill Location: ME

Instrument ID: B BFB Injection Date: 5/22/06

GC Column: RTX-502.2 ID: 0.25(mm) BFB Injection Time: 1030

Heated Purge: (Y/N) N

	IS1 AREA #	RT #	IS2 AREA #	RT #	IS3 AREA #	RT #	IS4 AREA #	RT #
12 HOUR STD	253263	7.00	428236	8.48	378997	13.02	180603	16.73
UPPER LIMIT	506526	7.42	856472	8.99	757994	13.80	361206	17.74
LOWER LIMIT	126632	6.58	214118	7.97	189499	12.23	90302	15.73
SAMPLE NO.								
01 0.2 PPB	211683	6.99	406327	8.48	364197	13.02	183801	16.74
02 1 PPB	250423	7.00	437021	8.49	379373	13.02	185550	16.74
03 2 PPB	257975	7.00	426815	8.48	375633	13.02	188191	16.74
04 5 PPB	251193	7.00	421056	8.48	376549	13.01	181024	16.74
05 20 PPB	253263	7.00	428236	8.48	378997	13.02	180603	16.73
06 50 PPB	239425	7.00	405568	8.49	362863	13.01	178724	16.74
07 100 PPB	221068	6.98	392598	8.47	366401	13.02	173234	16.74
08 ZZZZZ	264345	7.00	455555	8.49	411163	13.01	203347	16.73
09 ZZZZZ	255058	6.99	455989	8.48	411541	13.01	194605	16.74
10 ZZZZZ	271748	7.00	462940	8.48	412275	13.02	199284	16.74
11 ZZZZZ	272393	7.00	445229	8.49	385070	13.01	188312	16.73
12 ZZZZZ	260976	7.00	444257	8.48	394919	13.01	189595	16.73
13 ZZZZZ	263679	6.99	470007	8.48	411053	13.01	201605	16.74
14 ZZZZZ	274534	7.00	479456	8.48	421067	13.01	207298	16.74
15 ZZZZZ	276510	7.01	450004	8.49	405160	13.01	201256	16.73
16 ZZZZZ	280679	7.01	460882	8.48	410600	13.01	200215	16.74
17 ZZZZZ	273174	7.00	464748	8.48	409487	13.0045	202753	16.73
18 ZZZZZ	259604	7.00	446026	8.48	398451	13.0032	197966	16.73
19 ZZZZZ	270397	6.99	443005	8.48	396758	13.0097	194596	16.73
20 ZZZZZ	271294	6.99	457033	8.47	390557	13.0094	194644	16.73
21 ZZZZZ	274017	6.99	467123	8.48	405939	13.0056	197390	16.72
22 ZZZZZ	279785	6.99	460076	8.48	412418	13.0092	196815	16.73
23 ZZZZZ	281167	6.99	458846	8.48	412092	13.0088	199238	16.73
24 ZZZZZ	280997	7.00	481782	8.47	410598	13.001	201420	16.73

IS1 = Pentafluorobenzene
IS2 = Fluorobenzene
IS3 = Chlorobenzer
IS4 = 1,4-Dichlorobenzene-d4

AREA UPPER LIMIT = +100% of internal standard area
AREA LOWER LIMIT = -50% of internal standard area
RT UPPER LIMIT = +0.06 minutes of internal standard RRT
RT LOWER LIMIT = -0.06 minutes of internal standard RRT

Column used to flag values outside QC limits with an asterisk
* Values outside of QC limits.

analytics environmental laboratory LLC

VOLATILE ORGAN
INTERNAL STANDARD AREA AND RT SUM

Project #: 218736.002 Quote #: SDG#: 56637

BFB File ID: B36407T.D Site ID: GP Old Town Mill Location: ME

Instrument ID: B BFB Injection Date: 5/26/06

GC Column: RTX-502.2 ID: 0.25(mm) BFB Injection Time: 0956

Heated Purge: (Y/N) N

		IS1 AREA #	RT #	IS2 AREA #	RT #	IS3 AREA #	RT #	IS4 AREA #	RT #
	12 HOUR STD	99516	7.11	154615	8.54	134329	12.84	75097	16.34
	UPPER LIMIT	199032	7.53	309230	9.05	268658	13.61	150194	17.32
	LOWER LIMIT	49758	6.68	77308	8.02	67165	12.07	37549	15.36
	SAMPLE NO.								
01	1 PPB	92853	7.11	147055	8.54	124955	12.84	68277	16.33
02	2 PPB	93127	7.11	147758	8.54	124425	12.84	69352	16.34
03	5 PPB	93722	7.11	148737	8.55	126711	12.84	70309	16.34
04	20 PPB	99516	7.11	154615	8.54	134329	12.84	75097	16.34
05	50 PPB	106866	7.10	162033	8.54	140743	12.84	79036	16.33
06	100 PPB	111566	7.10	167116	8.54	147539	12.84	84904	16.34
07	ZZZZZ	106352	7.11	165177	8.54	138870	12.84	76053	16.33
08	ZZZZZ	108214	7.11	168759	8.55	145097	12.84	80027	16.34
09	ZZZZZ	109351	7.11	168831	8.55	144011	12.84	80704	16.34
10	ZZZZZ	104842	7.10	164284	8.54	138882	12.84	74778	16.33
11	ZZZZZ	101405	7.11	158267	8.54	135500	12.84	73232	16.34
12	ZZZZZ	100281	7.11	156659	8.54	134362	12.84	71760	16.34
13	ZZZZZ	97684	7.11	153838	8.54	130332	12.84	71195	16.34
14	ZZZZZ	96101	7.11	152756	8.54	130118	12.84	69884	16.34
15	ZZZZZ	96241	7.11	153427	8.55	130203	12.84	69766	16.34
16	ZZZZZ	96525	7.11	151513	8.55	130404	12.84	70542	16.34
17	ZZZZZ	100839	7.10	158284	8.54	135944	12.8415	76336	16.34
18	ZZZZZ	105080	7.11	163395	8.55	137948	12.8356	77703	16.34
19	ZZZZZ	98694	7.11	155601	8.55	132660	12.8358	71410	16.34
20	ZZZZZ	96073	7.11	152806	8.55	130650	12.8349	70826	16.34
21	ZZZZZ	96567	7.11	151389	8.55	131600	12.8354	70277	16.34
22	ZZZZZ	94429	7.11	150861	8.54	128965	12.8399	70680	16.34
23	ZZZZZ	96206	7.11	151676	8.55	129552	12.8373	70252	16.34
24									

IS1 = Pentafluorobenzene
IS2 = Fluorobenzene
IS3 = Chlorobenzer
IS4 = 1,4-Dichlorobenzene-d4

AREA UPPER LIMIT = +100% of internal standard area
AREA LOWER LIMIT = -50% of internal standard area
RT UPPER LIMIT = +0.06 minutes of internal standard RRT
RT LOWER LIMIT = -0.06 minutes of internal standard RRT

Column used to flag values outside QC limits with an asterisk
* Values outside of QC limits.

VOLATILE ORGAN
INTERNAL STANDARD AREA AND RT SUM

Project #: 218736.002 Quote #: SDG#: 56637

BFB File ID: B36458T.D Site ID: GP Old Town Mill Location: ME

Instrument ID: B BFB Injection Date: 5/30/06

GC Column: RTX-502.2 ID: 0.25(mm) BFB Injection Time: 0858

Heated Purge: (Y/N) N

		IS1 AREA #	RT #	IS2 AREA #	RT #	IS3 AREA #	RT #	IS4 AREA #	RT #
	12 HOUR STD	122493	7.11	189964	8.54	163115	12.84	92824	16.34
	UPPER LIMIT	244986	7.53	379928	9.05	326230	13.61	185648	17.32
	LOWER LIMIT	61247	6.68	94982	8.02	81558	12.07	46412	15.36
	SAMPLE NO.								
01	20 PPB	122493	7.11	189964	8.54	163115	12.84	92824	16.34
02	L805306B	121019	7.10	186897	8.54	163340	12.84	92305	16.33
03	ZZZZZ	120246	7.11	186387	8.55	163504	12.84	88726	16.34
04	B805306B	116009	7.11	181821	8.54	156910	12.84	84598	16.34
05	56637-29	112537	7.10	177637	8.54	154199	12.84	83327	16.34
06	ZZZZZ	111249	7.11	174561	8.54	150884	12.84	82155	16.34
07	ZZZZZ	110712	7.11	173864	8.55	149992	12.83	80212	16.34
08	ZZZZZ	107467	7.10	170873	8.54	150783	12.84	81243	16.33
09	ZZZZZ	107533	7.10	172383	8.54	148971	12.84	80687	16.33
10	ZZZZZ	107411	7.11	172104	8.54	150044	12.84	81925	16.34
11	56637-1	107359	7.11	169235	8.54	148789	12.84	80423	16.34
12	56637-2	104771	7.11	165802	8.54	144789	12.84	78561	16.34
13	56637-2,MS	113834	7.10	176676	8.54	153905	12.84	86113	16.34
14	56637-2,MSD	117505	7.10	181892	8.54	158258	12.84	87927	16.34
15	56637-3	109129	7.11	172196	8.55	148359	12.84	80408	16.34
16	56637-5	103990	7.11	162073	8.54	142928	12.84	78327	16.33
17	56637-20	106374	7.10	167995	8.54	146452	12.8342	78064	16.33
18	56637-21	105495	7.11	166550	8.55	146502	12.8355	79705	16.34
19	56637-24	105640	7.11	167193	8.54	144657	12.8363	78352	16.34
20	56637-25	106111	7.10	168811	8.54	144245	12.839	78898	16.34
21	56637-26	106127	7.11	165999	8.54	144593	12.8352	78599	16.34
22	56637-27	102924	7.10	165600	8.54	143514	12.8411	77644	16.33
23	56637-28	102774	7.10	161390	8.54	141639	12.8406	77351	16.33
24									

IS1 = Pentafluorobenzene
IS2 = Fluorobenzene
IS3 = Chlorobenzer
IS4 = 1,4-Dichlorobenzene-d4

AREA UPPER LIMIT = +30% of internal standard area
AREA LOWER LIMIT = -30% of internal standard area
RT UPPER LIMIT = +0.06 minutes of internal standard RRT
RT LOWER LIMIT = -0.06 minutes of internal standard RRT

Column used to flag values outside QC limits with an asterisk
* Values outside of QC limits.

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VOLATILE ORGAN
INTERNAL STANDARD AREA AND RT SUM

Project #: Quote #: SDG#:

BFB File ID: C59117T.D Site ID: Location:

Instrument ID: B BFB Injection Date: 5/30/06

GC Column: RTX-502.2 ID: 0.25(mm) BFB Injection Time: 0818

Heated Purge: (Y/N) N

		IS1 AREA #	RT #	IS2 AREA #	RT #	IS3 AREA #	RT #	IS4 AREA #	RT #
	12 HOUR STD	269108	6.94	457043	8.44	414522	12.97	199606	16.70
	UPPER LIMIT	538216	7.36	914086	8.95	829044	13.75	399212	17.70
	LOWER LIMIT	134554	6.53	228522	7.94	207261	12.19	99803	15.69
	SAMPLE NO.								
01	20 PPB	269108	6.94	457043	8.44	414522	12.97	199606	16.70
02	LS05306C	246240	6.94	421087	8.42	382230	12.97	197650	16.70
03	LS05306C2	255970	6.94	415541	8.43	380839	12.97	194940	16.70
04	ZZZZZ	281281	6.95	463185	8.44	413357	12.97	212342	16.69
05	MB05306C	262375	6.94	424504	8.43	395443	12.97	209356	16.70
06	56637-4	246605	6.94	432431	8.43	393378	12.97	207898	16.69
07	56637-14	269261	6.94	428970	8.43	382771	12.96	188223	16.69
08	56637-15	271078	6.94	451170	8.43	398927	12.97	196697	16.69
09	56637-16	268129	6.95	447163	8.43	393056	12.97	186576	16.70
10	ZZZZZ	237608	6.94	409003	8.43	360621	12.96	174402	16.69
11	56637-19	245317	6.95	417890	8.43	377476	12.97	187544	16.70
12	56637-9	261302	6.94	435022	8.43	393758	12.97	202186	16.69
13	56637-10	284023	6.96	459482	8.44	423240	12.98	214795	16.71
14	56637-11	283350	6.94	450283	8.43	411312	12.97	215147	16.70
15	ZZZZZ	273122	6.94	440186	8.43	396485	12.96	203256	16.69
16	ZZZZZ	264512	6.94	472051	8.43	413827	12.97	200453	16.70
17	56637-14, RR	282543	6.95	450336	8.43	417086	12.9648	206258	16.69
18	56637-17, RR	281807	6.94	443751	8.43	401051	12.9698	211660	16.70
19	ZZZZZ	281383	6.95	443955	8.43	399447	12.9705	204521	16.69
20	56637-18, RR,, 10ML FV	271033	6.95	433837	8.44	383990	12.9661	191274	16.69
21	ZZZZZ	278023	6.95	445774	8.43	405359	12.9606	207841	16.69
22	ZZZZZ	276783	6.95	465765	8.44	413905	12.9624	208205	16.69
23									
24									

IS1 = Pentafluorobenzene
IS2 = Fluorobenzene
IS3 = Chlorobenzer
IS4 = 1,4-Dichlorobenzene-d4

AREA UPPER LIMIT = +100% of internal standard area
AREA LOWER LIMIT = -50% of internal standard area
RT UPPER LIMIT = +0.06 minutes of internal standard RRT
RT LOWER LIMIT = -0.06 minutes of internal standard RRT

Column used to flag values outside QC limits with an asterisk
* Values outside of QC limits.

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EXTRACTION
LOGBOOK PAGES

Transfer Date: 05/30/06

Initials: AK

EXTRACT TRANSFER SHEET

Semi-Volatile Compounds:

VKM 5/30/06

JBH 05.30.06

Sample	Analysis	Extraction Date	Matrix	Initial Volume(mL)	Initial Weight(g)	Final Volume(mL)	% Solids	Extraction Comments
56637-1	8270W	05/30/06	Aqueous	1000		1.0		Emulsion
56637-2	8270W	05/30/06	Aqueous	1000		1.0		Emulsion
56637-3	8270W	05/30/06	Aqueous	1000		1.0		Emulsion
56637-20	8270W	05/30/06	Aqueous	1000		1.0		Emulsion
56637-21	8270W	05/30/06	Aqueous	1000		1.0		Emulsion

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Base Nuetral Acid Extractables Aqueous Preparation Logbook

Method References: SW-846 3510 Liquid/Liquid extraction; EPA 625

Date: 5/30/06
Time: 6:35 Initials: AK

QC Batch ID: Q-P-052606-1-AK
Prep. Batch ID P-P-053006-1-AK

Sample #	Initial Vol (mL)	mL Surrogate	mL Spike	1st ext. pH/Verify	Ext Vol Recov.	2nd ext. pH/verify	Total vol. Rec.	Final Vol $MeCl_2$	Cook Down Date
B05306AW	1000	0.5 2m	—	12.5 12.5	>150	1.5 1.5	>150	1mL	5/30/06 PM
L05306AW	1000	0.5	1.0 2m	12.5 12.5	>150	1.5 1.5	>150		
LD05306AW	1000	0.5	1.0 2m	12.5 12.5	>150	1.5 1.5	>150		
56636-1		0.5							PM 5/31
X 56637-2 *	1000	0.5	—	12.5 12.5	>150	1.5 1.5	>150	1mL	
X -3 *	1000	0.5	—	12.5 12.5	>150	1.5 1.5	>150	1	
X -20 *	1000	0.5	—	12.5 12.5	>150	1.5 1.5	>150		
-21 *	1000	0.5	—	12.5 12.5	>150	1.5 1.5			
X -1 *	1000	0.5	—	12.5 12.5	150	1.5 1.5	>150	1mL	5/30/06 AK

PM 6/2/06

* emulsions
X - 56637-1, -2, -3, -20 on 2nd Acid Shake, shaker table got turned up accidentally to 70. Some Sample (~5-10mL) was spilled.

MeCl2 Lot Number: 46066
1:1 H_2SO_4 Lot Number: M5382
NaOH Lot Number: 5205
Na_2SO_4 Lot Number: 055643

	Lot #	Exp Date
(1) ABN Surrogate:	56860	11/30/06
(2) ABN Spike (40 ug/mL):	56803	10/20/06
(3) Benzidines Spike (80 ug/mL):	56852	11/18/06
PAH Surrogate (5.0 ug/mL):		
PAH MS (5.0 ug/mL):		
TCLP MS (80 ug/mL):		

✓ DH 05.30.06

✓ BH 05-30-06

Transfer Date: 05/30/06

Initials: BH

EXTRACT TRANSFER SHEET

Semi-Volatile Compounds:

Sample	Analysis	Extraction Date	Matrix	Initial Volume(mL)	Initial Weight(g)	Final Volume(mL)	% Solids	Extraction Comments
56623-1	8270S	05/30/06	Solid		30.13	1.0	36.48	
56624-1	8270S	05/30/06	Solid		30.80	1.0	29.70	
56633-1	8270PAHS	05/30/06	Solid		30.70	1.0		
56633-2	8270PAHS	05/30/06	Solid		30.03	1.0		
56637-4	8270S	05/30/06	Solid		30.10	1.0		
56637-9	8270S	05/30/06	Solid		30.34	1.0		
56637-10	8270S	05/30/06	Solid		30.62	1.0		
56637-11	8270S	05/30/06	Solid		30.13	1.0		
56637-12	8270S	05/30/06	Solid		30.05	1.0		
56637-14	8270S	05/30/06	Solid		30.10	1.0		
56637-15	8270S	05/30/06	Solid		30.07	1.0		
56637-16	8270S	05/30/06	Solid		30.83	1.0		Excessive sodium sulfate required
56637-17	8270S	05/30/06	Solid		30.22	1.0		Excessive sodium sulfate required
56637-18	8270S	05/30/06	Solid		30.07	1.0		
56637-19	8270S	05/30/06	Solid		30.12	1.0		
56637-22	8270S	05/30/06	Solid		30.54	1.0		
56637-23	8270S	05/30/06	Solid		30.53	1.0		

✓ JJS 5-30-06

Base Nuetral Acid Extractables Solid Preparation Logbook

analytics environmental laboratory LLC

Circle Method Reference SW-846 3550B Sonication
SW-846 3545 Accelerated Solvent Extraction (ASE)

Date: 05·30·06 Time: 0805 QC Batch ID: Q-P-053006-1-BH
Initials: BH Prep. Batch ID P-P-053006-1-BH

Sample #	Matrix	Weight (g)	mL Surrogate ①	mL Spike ②/③	Final Vol $MeCl_2$	Cook Down Date	Vessel #
B05306AS	Na_2SO_4	30.98	0.5	—	1 mL	5·30·06	N/A
L05306AS	S. Sand	30.62	0.5	1.0/1.0		BH	
LD05306AS	↓	30.04	0.5	1.0/1.0			
56623-1	solid	30.13	0.5	—			
56624-1	soil	30.80	0.5	—			
56633-1	soil	30.70	0.5	—		05·30·06	
-2		30.03	0.5	—		BH	
56637-4		30.10	0.5	—			
-9		30.34	0.5	—			
-10		30.62	0.5	—			
-11		30.13	0.5	—			
-12		30.05	0.5	—			
-14		30.10	0.5	—			
-15		30.07	0.5	—			
-16 *		30.83	0.5	—			
-17 *		30.22	0.5	—			
-18		30.07	0.5	—			
-19		30.12	0.5	—			
-22		30.54	0.5	—			
-23		30.53 BH 05·30·06	0.5	—			
-23, ms		30.21	0.5	1.0/1.0			
-23, msd		30.20	0.5	1.0/1.0			

BH 05·30·06

$MeCl_2$ Lot Number: 46066
1:1 Acetone/MeCl2 Lot #: /
Silica Sand Lot Number: 074617
Na_2SO_4 Lot Number: 055643

	Lot #	Exp Date
① ABN Surrogate:	S6860	11·30·06
② ABN Spike (40 ug/mL):	S6803	10·20·06
③ Benzidines Spike (80 ug/mL):	S6832	11·18·06
PAH Surrogate (5.0 ug/mL):	/	/
PAH MS (5.0 ug/mL):	/	/

✓ JJ 5·30·06

* excessive sodium sulfate
[illegible]

Rev 1 01/16/04

analytics environmental laboratory LLC

SEMI-VOLATILE
DATA SUMMARIES

analytics environmental laboratory LLC

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800-929-9906

Mr. Karl Kasper
Woodard & Curran
41 Hutchins Drive
Portland ME 04102

June 6, 2006

CLIENT SAMPLE ID

Project Name: GP Old Town Mill

Project Number: 218736.002

Field Sample ID: LABQC

SAMPLE DATA

Lab Sample ID: B05306AW
Matrix: Aqueous
Percent Solid: N/A
Dilution Factor: 1.0
Collection Date: N/A
Lab Receipt Date: N/A
Extraction Date: 05/30/06
Analysis Date: 05/30/06

ANALYTICAL RESULTS SEMI-VOLATILE ORGANICS

ACID COMPOUND	Quantitation Limit µg/kg	Result µg/kg	ACID COMPOUND	Quantitation Limit µg/kg	Result µg/kg
2-Chlorophenol	5	U	Pentachlorophenol	10	U
4-Chloro-3-methylphenol	10	U	Phenol	5	U
2,4-Dichlorophenol	5	U	2,4,5-Trichlorophenol	5	U
2,4-Dimethylphenol	5	U	2,4,6-Trichlorophenol	5	U
2,4-dinitrophenol	5	U	Benzoic Acid	10	U
4,6-Dinitro-2-methylphenol	5	U	2-Methylphenol	5	U
2-Nitrophenol	5	U	3+4-Methylphenol	5	U
2,6-Dichlorophenol	5	U	Benzyl Alcohol	5	U
4-Nitrophenol	5	U	2,3,4,6-Tetrachlorophenol	5	U

Acid Surrogate Standard Recovery

2-Fluorophenol	51 %	d5-Phenol	37 %	2,4,6-Tribromophenol	85 %

BASE NEUTRAL COMPOUND	Quantitation Limit µg/kg	Result µg/kg	BASE NEUTRAL COMPOUND	Quantitation Limit µg/kg	Result µg/kg
1,2-Dichlorobenzene	2	U	Hexachlorobenzene	2	U
1,3-Dichlorobenzene	2	U	* Benzidine	20	U
1,4-Dichlorobenzene	2	U	3,3'-Dichlorobenzidine	20	U
2,4-Dinitrotoluene	2	U	Azobenzene	2	U
2,6-Dinitrotoluene	2	U	Bis(2-chloroethoxy)methane	2	U
Nitrobenzene	2	U	bis(2-chloroethyl) ether	2	U
Hexachlorobutadiene	2	U	bis(2-chloroisopropyl)ether	2	U
Dimethyl Phthalate	2	U	4-bromophenyl phenyl ether	2	U
Di-n-butyl phthalate	2	U	Butyl benzyl phthalate	2	U
di-n-octyl-phthalate	2	U	4-Chlorophenyl phenyl ether	2	U
Bis (2-ethylhexyl) phthalate	2	U	Diethyl Phthalate	2	U
1,2,4-Trichlorobenzene	2	U	Hexachlorocyclopentadiene	2	U

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis was conducted according to: Test Methods for Evaluating Solid Waste, SW-846 Method 8270C.

Authorized signature

8270/625 layout

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Mr. Karl Kasper
Woodard & Curran
41 Hutchins Drive
Portland ME 04102

June 6, 2006

CLIENT SAMPLE ID

Project Name: GP Old Town Mill

Project Number: 218736.002

Field Sample ID: LABQC

SAMPLE DATA

Lab Sample ID: B05306AW
Matrix: Aqueous
Percent Solid: N/A
Dilution Factor: 1.0
Collection Date: N/A
Lab Receipt Date: N/A
Extraction Date: 05/30/06
Analysis Date: 05/30/06

ANALYTICAL RESULTS SEMI-VOLATILE ORGANICS

BASE NEUTRAL COMPOUND	Quantitation Limit µg/kg	Result µg/kg	BASE NEUTRAL COMPOUND	Quantitation Limit µg/kg	Result µg/kg
Acenaphthene	2	U	N-nitrosodimethylamine	2	U
Acenaphthylene	2	U	N-nitroso-di-n-propylamine	2	U
Anthracene	2	U	n-nitrosodiphenylamine	2	U
Benzo[a]anthracene	2	U	Pyridine	2	U
Benzo[a] pyrene	2	U	2-Methylnaphthalene	2	U
Benzo[b] fluoranthene	2	U	2-Chloronaphthalene	2	U
Benzo[k] fluoranthene	2	U	Naphthalene	2	U
Benzo(g,h,i) perylene	2	U	Phenanthrene	2	U
Chrysene	2	U	Dibenzofuran	2	U
Dibenz [a,h] anthracene	2	U	Aniline	2	U
Fluoranthene	2	U	4-Chloroaniline	2	U
Fluorene	2	U	2-Nitroaniline	2	U
Indeno [1,2,3-cd] pyrene	2	U	3-Nitroaniline	2	U
Pyrene	2	U	4-Nitroaniline	2	U
Hexachloroethane	2	U	Carbazole	2	U
Isophorone	2	U			

Base Neutral Surrogate Standard Recovery

2-Fluorobiphenyl	75	%	d5-nitrobenzene	81	%	d14-p-terphenyl	84 %

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis was conducted according to: Test Methods for Evaluating Solid Waste, SW-846 Method 8270C.

COMMENTS: *Due to the reactive nature of this compound, the Benzidine quantitation limit is estimated.

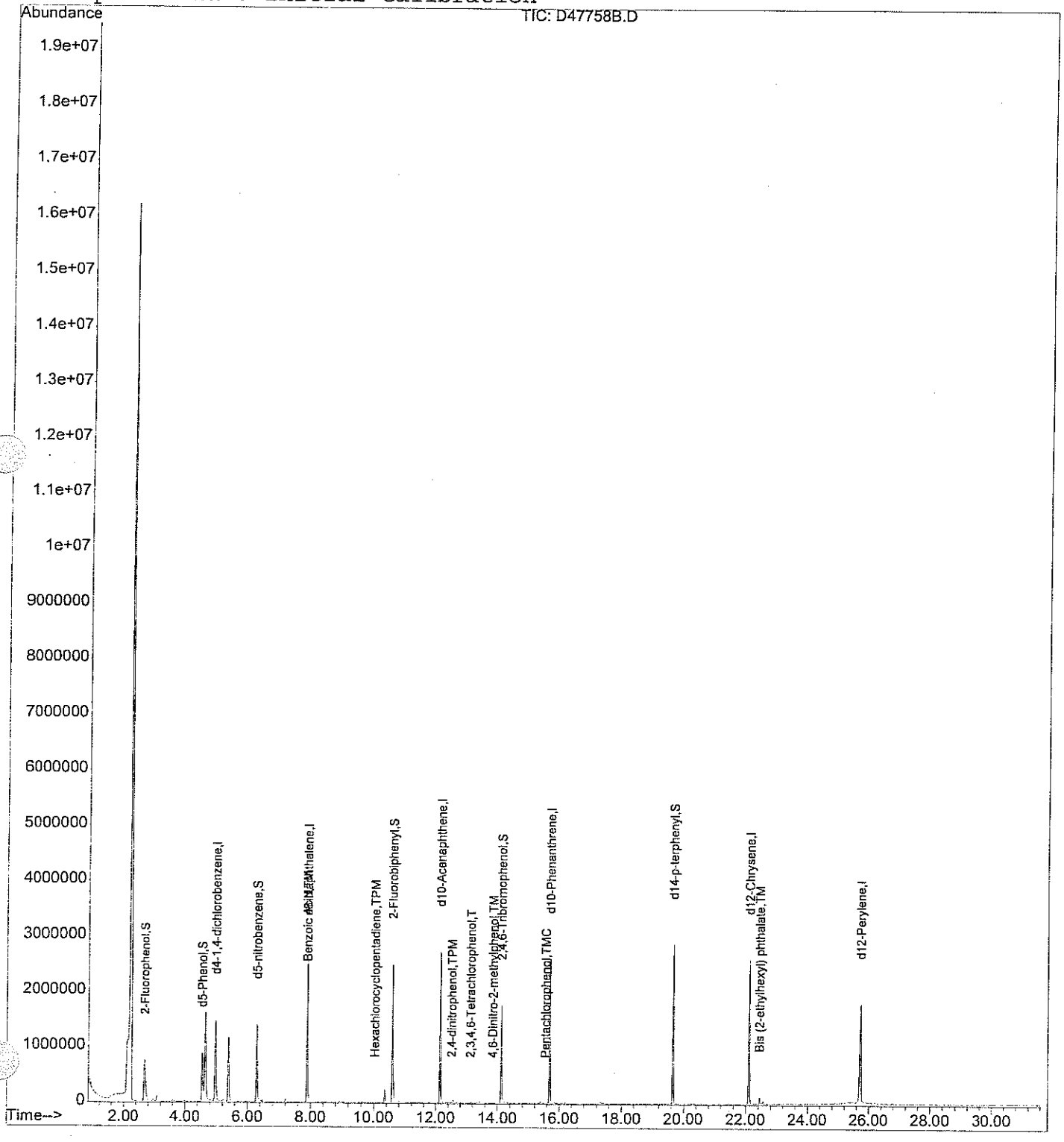
8270/625 layout

Authorized signature

Quantitation Report

Data File : D:\HPCHEM\1\DATA\053006-D\D47758B.D Vial: 4
Acq On : 30 May 2006 9:36 am Operator:
Sample : B05306AW Inst : D Instrum
Misc : Multiplr: 1.00
MS Integration Params: RTEINT.P
Quant Time: May 30 9:52 2006 Quant Results File: S05246.RES

Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Tue May 30 09:52:41 2006
Response via : Initial Calibration



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800-929-9906

Mr. Karl Kasper
Woodard & Curran
41 Hutchins Drive
Portland ME 04102

June 6, 2006

CLIENT SAMPLE ID

Project Name: GP Old Town Mill

Project Number: 218736.002

Field Sample ID: LABQC

SAMPLE DATA

Lab Sample ID: B05306AS
Matrix: Solid
Percent Solid: 100
Dilution Factor: 1.0
Collection Date: N/A
Lab Receipt Date: N/A
Extraction Date: 05/30/06
Analysis Date: 05/30/06

ANALYTICAL RESULTS SEMI-VOLATILE ORGANICS

ACID COMPOUND	Quantitation Limit µg/kg	Result µg/kg	ACID COMPOUND	Quantitation Limit µg/kg	Result µg/kg
2-Chlorophenol	250	U	Pentachlorophenol	250	U
4-Chloro-3-methylphenol	250	U	Phenol	250	U
2,4-Dichlorophenol	250	U	2,4,5-Trichlorophenol	250	U
2,4-Dimethylphenol	250	U	2,4,6-Trichlorophenol	250	U
2,4-dinitrophenol	250	U	Benzoic Acid	250	U
4,6-Dinitro-2-methylphenol	250	U	2-Methylphenol	250	U
2-Nitrophenol	250	U	3+4-Methylphenol	250	U
2,6-Dichlorophenol	250	U	Benzyl Alcohol	250	U
4-Nitrophenol	250	U	2,3,4,6-Tetrachlorophenol	250	U

Acid Surrogate Standard Recovery

2-Fluorophenol	68 %	d5-Phenol	77 %	2,4,6-Tribromophenol	77 %

BASE NEUTRAL COMPOUND	Quantitation Limit µg/kg	Result µg/kg	BASE NEUTRAL COMPOUND	Quantitation Limit µg/kg	Result µg/kg
1,2-Dichlorobenzene	250	U	Hexachlorobenzene	250	U
1,3-Dichlorobenzene	250	U	* Benzidine	250	U
1,4-Dichlorobenzene	250	U	3,3'-Dichlorobenzidine	250	U
2,4-Dinitrotoluene	250	U	Azobenzene	250	U
2,6-Dinitrotoluene	250	U	Bis(2-chloroethoxy)methane	250	U
Nitrobenzene	250	U	bis(2-chloroethyl) ether	250	U
Hexachlorobutadiene	250	U	bis(2-chloroisopropyl)ether	250	U
Dimethyl Phthalate	250	U	4-bromophenyl phenyl ether	250	U
Di-n-butyl phthalate	250	U	Butyl benzyl phthalate	250	U
di-n-octyl-phthalate	250	U	4-Chlorophenyl phenyl ether	250	U
Bis (2-ethylhexyl) phthalate	250	U	Diethyl Phthalate	250	U
1,2,4-Trichlorobenzene	250	U	Hexachlorocyclopentadiene	250	U

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis was conducted according to: Test Methods for Evaluating Solid Waste, SW-846 Method 8270C.

Authorized signature

8270/625 layout

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Mr. Karl Kasper
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June 6, 2006

CLIENT SAMPLE ID

Project Name: GP Old Town Mill

Project Number: 218736.002

Field Sample ID: LABQC

SAMPLE DATA

Lab Sample ID: B05306AS
Matrix: Solid
Percent Solid: 100
Dilution Factor: 1.0
Collection Date: N/A
Lab Receipt Date: N/A
Extraction Date: 05/30/06
Analysis Date: 05/30/06

ANALYTICAL RESULTS SEMI-VOLATILE ORGANICS

BASE NEUTRAL COMPOUND	Quantitation Limit µg/kg	Result µg/kg	BASE NEUTRAL COMPOUND	Quantitation Limit µg/kg	Result µg/kg
Acenaphthene	250	U	N-nitrosodimethylamine	250	U
Acenaphthylene	250	U	N-nitroso-di-n-propylamine	250	U
Anthracene	250	U	n-nitrosodiphenylamine	250	U
Benzo[a]anthracene	250	U	Pyridine	250	U
Benzo[a] pyrene	250	U	2-Methylnaphthalene	250	U
Benzo[b] fluoranthene	250	U	2-Chloronaphthalene	250	U
Benzo[k] fluoranthene	250	U	Naphthalene	250	U
Benzo(g,h,i) perylene	250	U	Phenanthrene	250	U
Chrysene	250	U	Dibenzofuran	250	U
Dibenz [a,h] anthracene	250	U	Aniline	250	U
Fluoranthene	250	U	4-Chloroaniline	250	U
Fluorene	250	U	2-Nitroaniline	250	U
Indeno [1,2,3-cd] pyrene	250	U	3-Nitroaniline	250	U
Pyrene	250	U	4-Nitroaniline	250	U
Hexachloroethane	250	U	Carbazole	250	U
Isophorone	250	U			

Base Neutral Surrogate Standard Recovery

2-Fluorobiphenyl	77	%	d5-nitrobenzene	72	%	d14-p-terphenyl	83 %

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis was conducted according to: Test Methods for Evaluating Solid Waste, SW-846 Method 8270C.

COMMENTS: *Due to the reactive nature of this compound, the Benzidine quantitation limit is estimated.
Results are expressed on a dry weight basis.

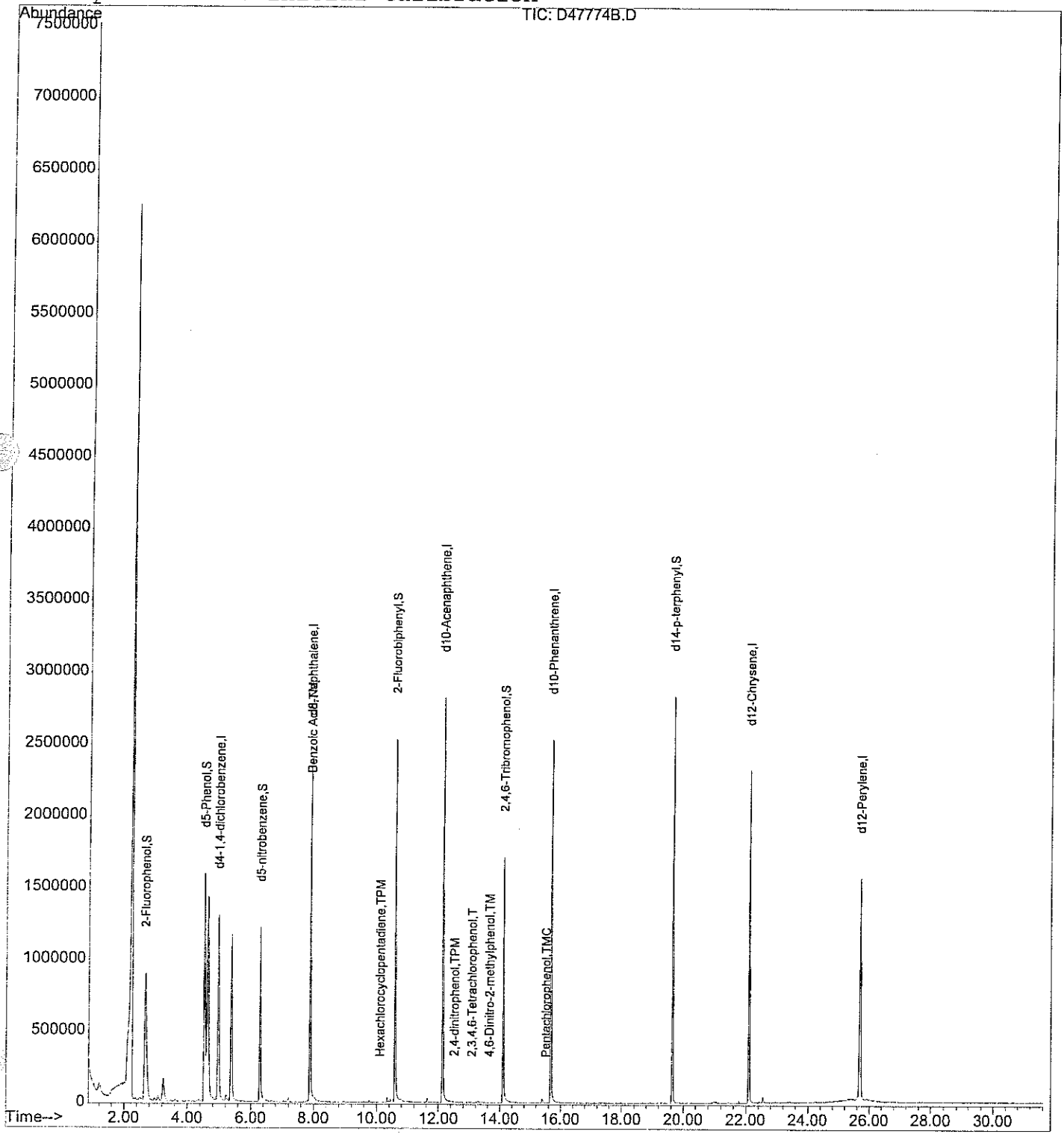
8270/625 layout

Authorized signature

Quantitation Report

Data File : D:\HPCHEM\1\DATA\053006-D\D47774B.D Vial: 4
Acq On : 30 May 2006 10:19 pm Operator:
Sample : B05306AS Inst : D Instrum
Misc : SOIL Multiplr: 1.00
MS Integration Params: RTEINT.P
Quant Time: May 31 8:20 2006 Quant Results File: S05246.RES

Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Tue May 30 20:31:56 2006
Response via : Initial Calibration



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41 Hutchins Drive
Portland ME 04102

June 6, 2006

CLIENT SAMPLE ID

Project Name: GP Old Town Mill

Project Number: 218736.002

Field Sample ID: LABQC

SAMPLE DATA

Lab Sample ID: B05306AS RR
Matrix: Solid
Percent Solid: 100
Dilution Factor: 1.0
Collection Date: N/A
Lab Receipt Date: N/A
Extraction Date: 05/30/06
Analysis Date: 05/31/06

ANALYTICAL RESULTS SEMI-VOLATILE ORGANICS

ACID COMPOUND	Quantitation Limit µg/kg	Result µg/kg	ACID COMPOUND	Quantitation Limit µg/kg	Result µg/kg
2-Chlorophenol	250	U	Pentachlorophenol	250	U
4-Chloro-3-methylphenol	250	U	Phenol	250	U
2,4-Dichlorophenol	250	U	2,4,5-Trichlorophenol	250	U
2,4-Dimethylphenol	250	U	2,4,6-Trichlorophenol	250	U
2,4-dinitrophenol	250	U	Benzoic Acid	250	U
4,6-Dinitro-2-methylphenol	250	U	2-Methylphenol	250	U
2-Nitrophenol	250	U	3+4-Methylphenol	250	U
2,6-Dichlorophenol	250	U	Benzyl Alcohol	250	U
4-Nitrophenol	250	U	2,3,4,6-Tetrachlorophenol	250	U

Acid Surrogate Standard Recovery

2-Fluorophenol	67 %	d5-Phenol	78 %	2,4,6-Tribromophenol	85 %

BASE NEUTRAL COMPOUND	Quantitation Limit µg/kg	Result µg/kg	BASE NEUTRAL COMPOUND	Quantitation Limit µg/kg	Result µg/kg
1,2-Dichlorobenzene	250	U	Hexachlorobenzene	250	U
1,3-Dichlorobenzene	250	U	* Benzidine	250	U
1,4-Dichlorobenzene	250	U	3,3'-Dichlorobenzidine	250	U
2,4-Dinitrotoluene	250	U	Azobenzene	250	U
2,6-Dinitrotoluene	250	U	Bis(2-chloroethoxy)methane	250	U
Nitrobenzene	250	U	bis(2-chloroethyl) ether	250	U
Hexachlorobutadiene	250	U	bis(2-chloroisopropyl)ether	250	U
Dimethyl Phthalate	250	U	4-bromophenyl phenyl ether	250	U
Di-n-butyl phthalate	250	U	Butyl benzyl phthalate	250	U
di-n-octyl-phthalate	250	U	4-Chlorophenyl phenyl ether	250	U
Bis (2-ethylhexyl) phthalate	250	U	Diethyl Phthalate	250	U
1,2,4-Trichlorobenzene	250	U	Hexachlorocyclopentadiene	250	U

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis was conducted according to: Test Methods for Evaluating Solid Waste, SW-846 Method 8270C.

Authorized signature Melissa Hall

8270/625 layout

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June 6, 2006

SAMPLE DATA

Lab Sample ID:	B05306AS	RR
Matrix:	Solid	
Percent Solid:	100	
Dilution Factor:	1.0	
Collection Date:	N/A	
Lab Receipt Date:	N/A	
Extraction Date:	05/30/06	
Analysis Date:	05/31/06	

CLIENT SAMPLE ID

Project Name:	GP Old Town Mill
Project Number:	218736.002
Field Sample ID:	LABQC

ANALYTICAL RESULTS SEMI-VOLATILE ORGANICS

BASE NEUTRAL COMPOUND	Quantitation Limit µg/kg	Result µg/kg	BASE NEUTRAL COMPOUND	Quantitation Limit µg/kg	Result µg/kg
Acenaphthene	250	U	N-nitrosodimethylamine	250	U
Acenaphthylene	250	U	N-nitroso-di-n-propylamine	250	U
Anthracene	250	U	n-nitrosodiphenylamine	250	U
Benzo[a]anthracene	250	U	Pyridine	250	U
Benzo[a] pyrene	250	U	2-Methylnaphthalene	250	U
Benzo[b] fluoranthene	250	U	2-Chloronaphthalene	250	U
Benzo[k] fluoranthene	250	U	Naphthalene	250	U
Benzo(g,h,i) perylene	250	U	Phenanthrene	250	U
Chrysene	250	U	Dibenzofuran	250	U
Dibenz [a,h] anthracene	250	U	Aniline	250	U
Fluoranthene	250	U	4-Chloroaniline	250	U
Fluorene	250	U	2-Nitroaniline	250	U
Indeno [1,2,3-cd] pyrene	250	U	3-Nitroaniline	250	U
Pyrene	250	U	4-Nitroaniline	250	U
Hexachloroethane	250	U	Carbazole	250	U
Isophorone	250	U			

Base Neutral Surrogate Standard Recovery

2-Fluorobiphenyl	77	%	d5-nitrobenzene	72	%	d14-p-terphenyl	86 %

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis was conducted according to: Test Methods for Evaluating Solid Waste, SW-846 Method 8270C.

COMMENTS: *Due to the reactive nature of this compound, the Benzidine quantitation limit is estimated.
Results are expressed on a dry weight basis.

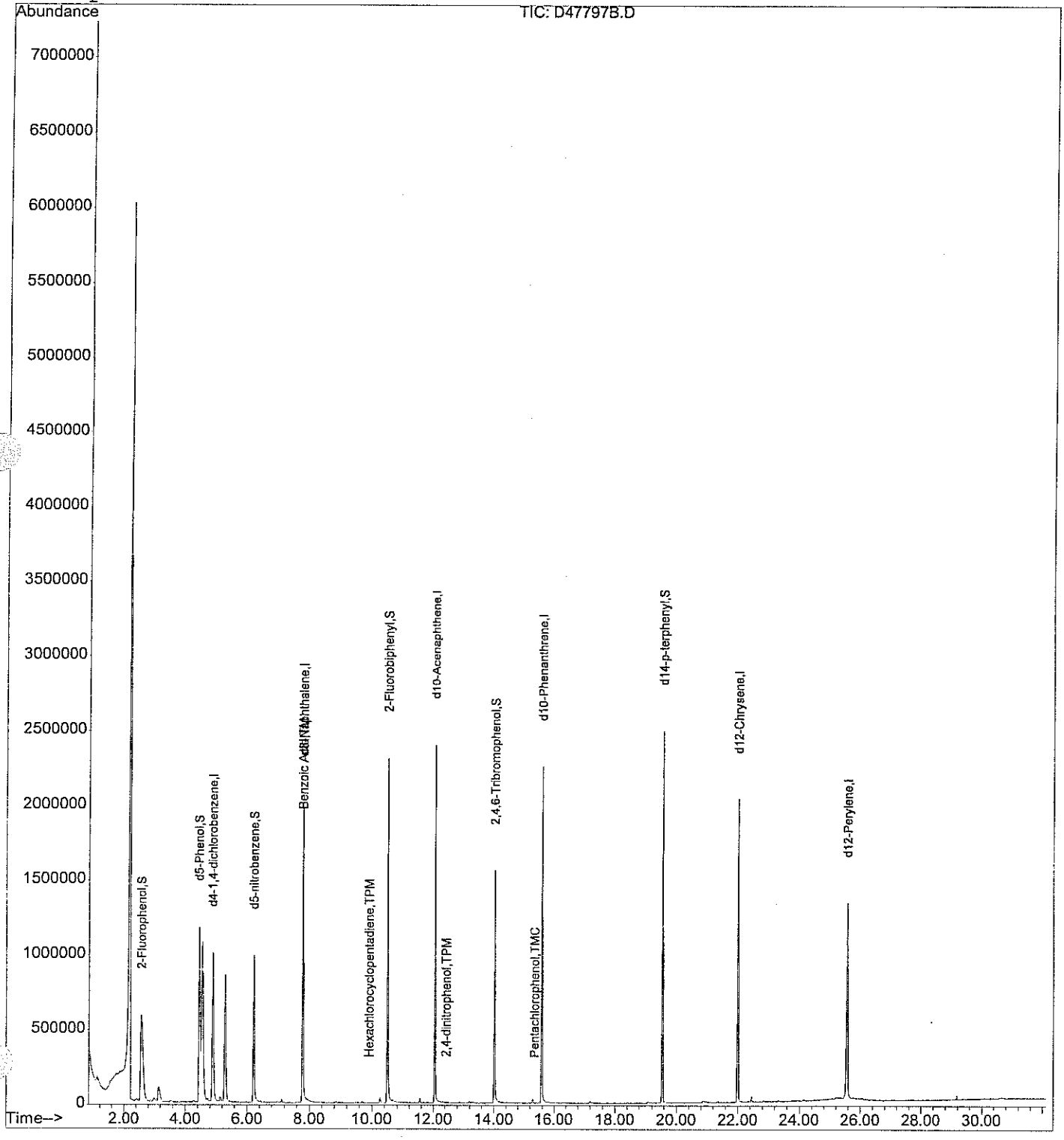
8270/625 layout

Authorized signature [signature]

Quantitation Report

Data File : D:\HPCHEM\1\DATA\053106-D\D47797B.D Vial: 4
Acq On : 31 May 2006 12:13 pm Operator:
Sample : B05306AS,RE Inst : D Instrum
Misc : SOIL Multiplr: 1.00
MS Integration Params: RTEINT.P
Quant Time: May 31 12:36 2006 Quant Results File: S05246.RES

Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Wed May 31 19:58:02 2006
Response via : Initial Calibration



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May 31, 2006

CLIENT SAMPLE ID

Project Name: GP Old Town Mill

Project Number: 218736.002

Field Sample ID: MW-5

SAMPLE DATA

Lab Sample ID: 56637-1
Matrix: Aqueous
Percent Solid: N/A
Dilution Factor: 1.0
Collection Date: 05/26/06
Lab Receipt Date: 05/27/06
Extraction Date: 05/30/06
Analysis Date: 05/30/06

ANALYTICAL RESULTS SEMI-VOLATILE ORGANICS

ACID COMPOUND	Quantitation Limit µg/L	Result µg/L	ACID COMPOUND	Quantitation Limit µg/L	Result µg/L
2-Chlorophenol	5	U	Pentachlorophenol	10	U
4-Chloro-3-methylphenol	10	U	Phenol	5	U
2,4-Dichlorophenol	5	U	2,4,5-Trichlorophenol	5	U
2,4-Dimethylphenol	5	U	2,4,6-Trichlorophenol	5	U
2,4-dinitrophenol	5	U	Benzoic Acid	10	U
4,6-Dinitro-2-methylphenol	5	U	2-Methylphenol	5	U
2-Nitrophenol	5	U	3+4-Methylphenol	5	U
2,6-Dichlorophenol	5	U	Benzyl Alcohol	5	U
4-Nitrophenol	5	U	2,3,4,6-Tetrachlorophenol	5	U

Acid Surrogate Standard Recovery

2-Fluorophenol	38 %	d5-Phenol	31 %	2,4,6-Tribromophenol	62* %

BASE NEUTRAL COMPOUND	Quantitation Limit µg/L	Result µg/L	BASE NEUTRAL COMPOUND	Quantitation Limit µg/L	Result µg/L
1,2-Dichlorobenzene	2	U	Hexachlorobenzene	2	U
1,3-Dichlorobenzene	2	U	* Benzidine	20	U
1,4-Dichlorobenzene	2	U	3,3'-Dichlorobenzidine	20	U
2,4-Dinitrotoluene	2	U	Azobenzene	2	U
2,6-Dinitrotoluene	2	U	Bis(2-chloroethoxy)methane	2	U
Nitrobenzene	2	U	bis(2-chloroethyl) ether	2	U
Hexachlorobutadiene	2	U	bis(2-chloroisopropyl)ether	2	U
Dimethyl Phthalate	2	U	4-bromophenyl phenyl ether	2	U
Di-n-butyl phthalate	2	U	Butyl benzyl phthalate	2	U
di-n-octyl-phthalate	2	U	4-Chlorophenyl phenyl ether	2	U
Bis (2-ethylhexyl) phthalate	2	4	Diethyl Phthalate	2	U
1,2,4-Trichlorobenzene	2	U	Hexachlorocyclopentadiene	2	U

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis was conducted according to: Test Methods for Evaluating Solid Waste, SW-846 Method 8270C.

Authorized signature Melena Tull

8270/625 layout

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May 31, 2006

CLIENT SAMPLE ID

Project Name: GP Old Town Mill

Project Number: 218736.002

Field Sample ID: MW-5

SAMPLE DATA

Lab Sample ID: 56637-1
Matrix: Aqueous
Percent Solid: N/A
Dilution Factor: 1.0
Collection Date: 05/26/06
Lab Receipt Date: 05/27/06
Extraction Date: 05/30/06
Analysis Date: 05/30/06

ANALYTICAL RESULTS SEMI-VOLATILE ORGANICS

BASE NEUTRAL COMPOUND	Quantitation Limit µg/L	Result µg/L	BASE NEUTRAL COMPOUND	Quantitation Limit µg/L	Result µg/L
Acenaphthene	2	U	N-nitrosodimethylamine	2	U
Acenaphthylene	2	U	N-nitroso-di-n-propylamine	2	U
Anthracene	2	U	n-nitrosodiphenylamine	2	U
Benzo[a]anthracene	2	**2**	Pyridine	2	U
Benzo[a] pyrene	2	**1 J**	2-Methylnaphthalene	2	U
Benzo[b] fluoranthene	2	**1 J**	2-Chloronaphthalene	2	U
Benzo[k] fluoranthene	2	**1 J**	Naphthalene	2	U
Benzo(g,h,i) perylene	2	U	Phenanthrene	2	**2**
Chrysene	2	**1 J**	Dibenzofuran	2	U
Dibenz [a,h] anthracene	2	U	Aniline	2	U
Fluoranthene	2	**3**	4-Chloroaniline	2	U
Fluorene	2	U	2-Nitroaniline	2	U
Indeno [1,2,3-cd] pyrene	2	U	3-Nitroaniline	2	U
Pyrene	2	**3**	4-Nitroaniline	2	U
Hexachloroethane	2	U	Carbazole	2	U
Isophorone	2	U			

Base Neutral Surrogate Standard Recovery

2-Fluorobiphenyl	75	%	d5-nitrobenzene	77	%	d14-p-terphenyl	48 %

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis was conducted according to: Test Methods for Evaluating Solid Waste, SW-846 Method 8270C.

COMMENTS: *Due to the reactive nature of this compound, the Benzidine quantitation limit is estimated.
*Surrogate recovery outside laboratory acceptance criteria. Sample was reanalyzed with similar results.

8270/625 layout

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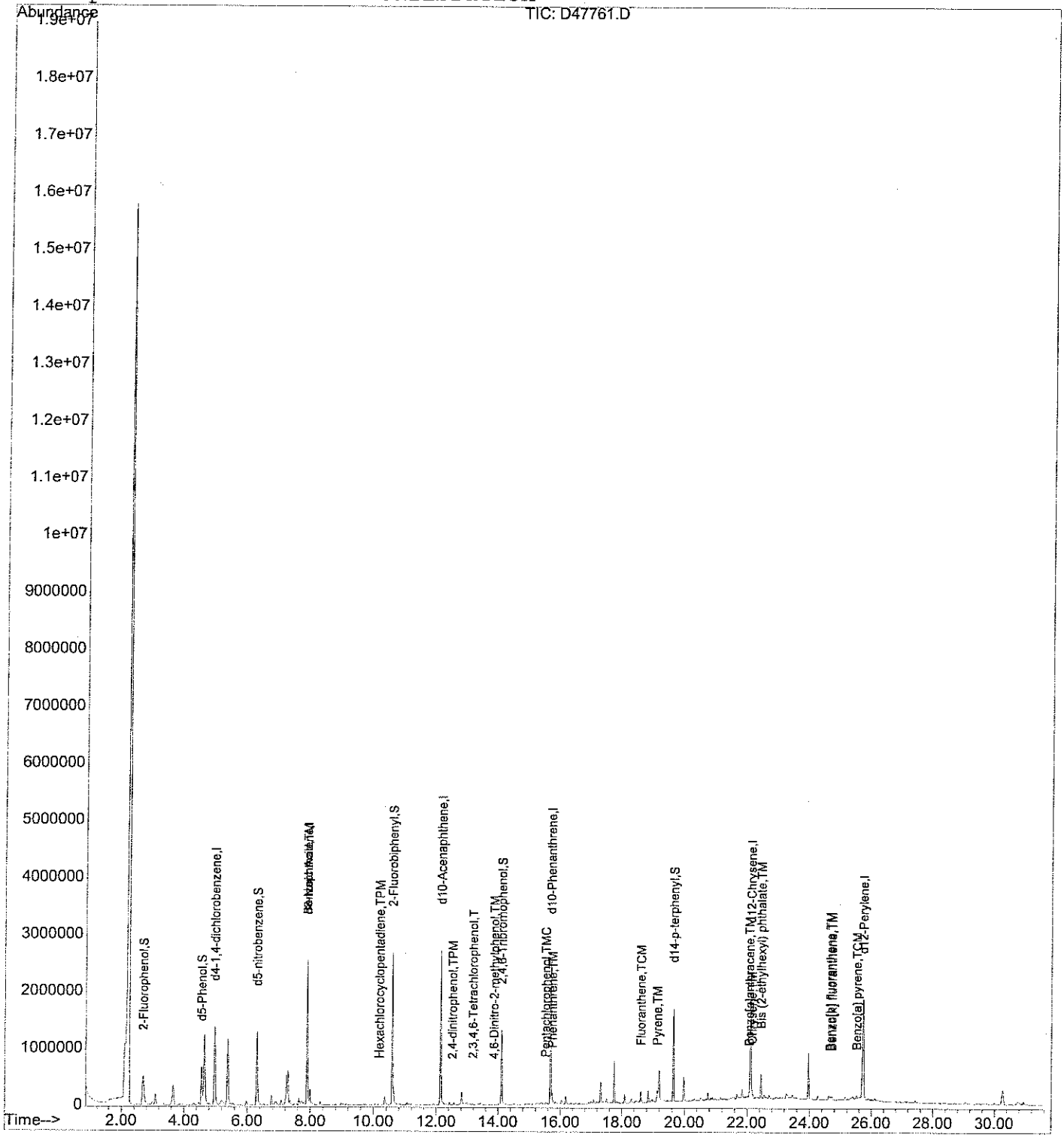
Quantitation Report

Data File : D:\HPCHEM\1\DATA\053006-D\D47761.D Vial: 7
Acq On : 30 May 2006 10:56 am Operator:
Sample : 56637-1 Inst : D Instrum
Misc : Multiplr: 1.00
MS Integration Params: RTEINT.P
Quant Time: May 30 11:32 2006 Quant Results File: S05246.RES

Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Tue May 30 09:52:41 2006
Response via : Initial Calibration



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May 31, 2006

CLIENT SAMPLE ID

Project Name: GP Old Town Mill

Project Number: 218736.002

Field Sample ID: MW-4

SAMPLE DATA

Lab Sample ID: 56637-2
Matrix: Aqueous
Percent Solid: N/A
Dilution Factor: 1.0
Collection Date: 05/26/06
Lab Receipt Date: 05/27/06
Extraction Date: 05/30/06
Analysis Date: 05/30/06

ANALYTICAL RESULTS SEMI-VOLATILE ORGANICS

ACID COMPOUND	Quantitation Limit µg/L	Result µg/L	ACID COMPOUND	Quantitation Limit µg/L	Result µg/L
2-Chlorophenol	5	U	Pentachlorophenol	10	U
4-Chloro-3-methylphenol	10	U	Phenol	5	U
2,4-Dichlorophenol	5	U	2,4,5-Trichlorophenol	5	U
2,4-Dimethylphenol	5	U	2,4,6-Trichlorophenol	5	U
2,4-dinitrophenol	5	U	Benzoic Acid	10	U
4,6-Dinitro-2-methylphenol	5	U	2-Methylphenol	5	U
2-Nitrophenol	5	U	3+4-Methylphenol	5	U
2,6-Dichlorophenol	5	U	Benzyl Alcohol	5	U
4-Nitrophenol	5	U	2,3,4,6-Tetrachlorophenol	5	U

Acid Surrogate Standard Recovery

2-Fluorophenol	2 * %	d5-Phenol	0 * %	2,4,6-Tribromophenol	8* %

BASE NEUTRAL COMPOUND	Quantitation Limit µg/L	Result µg/L	BASE NEUTRAL COMPOUND	Quantitation Limit µg/L	Result µg/L
1,2-Dichlorobenzene	2	U	Hexachlorobenzene	2	U
1,3-Dichlorobenzene	2	U	* Benzidine	20	U
1,4-Dichlorobenzene	2	U	3,3'-Dichlorobenzidine	20	U
2,4-Dinitrotoluene	2	U	Azobenzene	2	U
2,6-Dinitrotoluene	2	U	Bis(2-chloroethoxy)methane	2	U
Nitrobenzene	2	U	bis(2-chloroethyl) ether	2	U
Hexachlorobutadiene	2	U	bis(2-chloroisopropyl)ether	2	U
Dimethyl Phthalate	2	U	4-bromophenyl phenyl ether	2	U
Di-n-butyl phthalate	2	U	Butyl benzyl phthalate	2	U
di-n-octyl-phthalate	2	U	4-Chlorophenyl phenyl ether	2	U
Bis (2-ethylhexyl) phthalate	2	U	Diethyl Phthalate	2	U
1,2,4-Trichlorobenzene	2	U	Hexachlorocyclopentadiene	2	U

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis was conducted according to: Test Methods for Evaluating Solid Waste, SW-846 Method 8270C.

Authorized signature

8270/625 layout

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May 31, 2006

CLIENT SAMPLE ID

Project Name: GP Old Town Mill

Project Number: 218736.002

Field Sample ID: MW-4

SAMPLE DATA

Lab Sample ID: 56637-2
Matrix: Aqueous
Percent Solid: N/A
Dilution Factor: 1.0
Collection Date: 05/26/06
Lab Receipt Date: 05/27/06
Extraction Date: 05/30/06
Analysis Date: 05/30/06

ANALYTICAL RESULTS SEMI-VOLATILE ORGANICS

BASE NEUTRAL COMPOUND	Quantitation Limit µg/L	Result µg/L	BASE NEUTRAL COMPOUND	Quantitation Limit µg/L	Result µg/L
Acenaphthene	2	U	N-nitrosodimethylamine	2	U
Acenaphthylene	2	U	N-nitroso-di-n-propylamine	2	U
Anthracene	2	U	n-nitrosodiphenylamine	2	U
Benzo[a]anthracene	2	U	Pyridine	2	U
Benzo[a] pyrene	2	U	2-Methylnaphthalene	2	U
Benzo[b] fluoranthene	2	U	2-Chloronaphthalene	2	U
Benzo[k] fluoranthene	2	U	Naphthalene	2	U
Benzo(g,h,i) perylene	2	U	Phenanthrene	2	U
Chrysene	2	U	Dibenzofuran	2	U
Dibenz [a,h] anthracene	2	U	Aniline	2	U
Fluoranthene	2	U	4-Chloroaniline	2	U
Fluorene	2	U	2-Nitroaniline	2	U
Indeno [1,2,3-cd] pyrene	2	U	3-Nitroaniline	2	U
Pyrene	2	U	4-Nitroaniline	2	U
Hexachloroethane	2	U	Carbazole	2	U
Isophorone	2	U			

Base Neutral Surrogate Standard Recovery

2-Fluorobiphenyl	74	%	d5-nitrobenzene	76	%	d14-p-terphenyl	55 %

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis was conducted according to: Test Methods for Evaluating Solid Waste, SW-846 Method 8270C.

COMMENTS: *Due to the reactive nature of this compound, the Benzidine quantitation limit is estimated.
*Surrogate recovery outside laboratory acceptance criteria. Sample was reanalyzed with similar results.

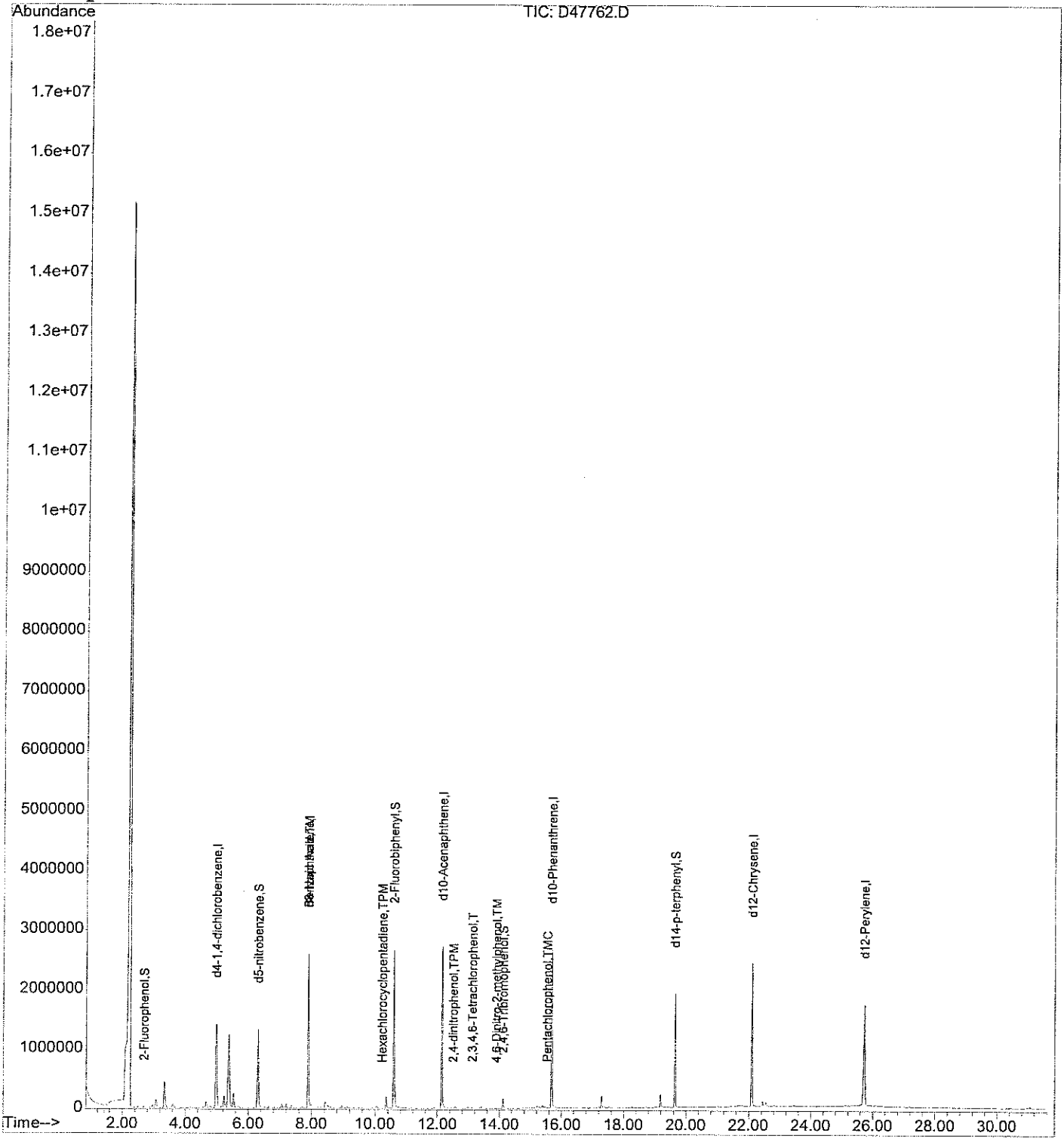
8270/625 layout

Authorized signature

Quantitation Report

Data File : D:\HPCHEM\1\DATA\053006-D\D47762.D Vial: 8
Acq On : 30 May 2006 12:02 pm Operator:
Sample : 56637-2 Inst : D Instrum
Misc : Multiplr: 1.00
MS Integration Params: RTEINT.P
Quant Time: May 30 12:34 2006 Quant Results File: S05246.RES

Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Tue May 30 09:52:41 2006
Response via : Initial Calibration



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May 31, 2006

CLIENT SAMPLE ID

Project Name: GP Old Town Mill

Project Number: 218736.002

Field Sample ID: SB-37

SAMPLE DATA

Lab Sample ID: 56637-3
Matrix: Aqueous
Percent Solid: N/A
Dilution Factor: 1.0
Collection Date: 05/26/06
Lab Receipt Date: 05/27/06
Extraction Date: 05/30/06
Analysis Date: 05/30/06

ANALYTICAL RESULTS SEMI-VOLATILE ORGANICS

ACID COMPOUND	Quantitation Limit µg/L	Result µg/L	ACID COMPOUND	Quantitation Limit µg/L	Result µg/L
2-Chlorophenol	5	U	Pentachlorophenol	10	U
4-Chloro-3-methylphenol	10	U	Phenol	5	U
2,4-Dichlorophenol	5	U	2,4,5-Trichlorophenol	5	U
2,4-Dimethylphenol	5	U	2,4,6-Trichlorophenol	5	U
2,4-dinitrophenol	5	U	Benzoic Acid	10	U
4,6-Dinitro-2-methylphenol	5	U	2-Methylphenol	5	U
2-Nitrophenol	5	U	3+4-Methylphenol	5	U
2,6-Dichlorophenol	5	U	Benzyl Alcohol	5	U
4-Nitrophenol	5	U	2,3,4,6-Tetrachlorophenol	5	U

Acid Surrogate Standard Recovery

2-Fluorophenol	40 %	d5-Phenol	29 %	2,4,6-Tribromophenol	82 %

BASE NEUTRAL COMPOUND	Quantitation Limit µg/L	Result µg/L	BASE NEUTRAL COMPOUND	Quantitation Limit µg/L	Result µg/L
1,2-Dichlorobenzene	2	U	Hexachlorobenzene	2	U
1,3-Dichlorobenzene	2	U	* Benzidine	20	U
1,4-Dichlorobenzene	2	U	3,3'-Dichlorobenzidine	20	U
2,4-Dinitrotoluene	2	U	Azobenzene	2	U
2,6-Dinitrotoluene	2	U	Bis(2-chloroethoxy)methane	2	U
Nitrobenzene	2	U	bis(2-chloroethyl) ether	2	U
Hexachlorobutadiene	2	U	bis(2-chloroisopropyl)ether	2	U
Dimethyl Phthalate	2	U	4-bromophenyl phenyl ether	2	U
Di-n-butyl phthalate	2	U	Butyl benzyl phthalate	2	U
di-n-octyl-phthalate	2	U	4-Chlorophenyl phenyl ether	2	U
Bis (2-ethylhexyl) phthalate	2	3	Diethyl Phthalate	2	U
1,2,4-Trichlorobenzene	2	U	Hexachlorocyclopentadiene	2	U

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis was conducted according to: Test Methods for Evaluating Solid Waste, SW-846 Method 8270C.

Authorized signature Milena Tull

8270/525 layout

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May 31, 2006

CLIENT SAMPLE ID

Project Name: GP Old Town Mill

Project Number: 218736.002

Field Sample ID: SB-37

SAMPLE DATA

Lab Sample ID: 56637-3
Matrix: Aqueous
Percent Solid: N/A
Dilution Factor: 1.0
Collection Date: 05/26/06
Lab Receipt Date: 05/27/06
Extraction Date: 05/30/06
Analysis Date: 05/30/06

ANALYTICAL RESULTS SEMI-VOLATILE ORGANICS

BASE NEUTRAL COMPOUND	Quantitation Limit µg/L	Result µg/L	BASE NEUTRAL COMPOUND	Quantitation Limit µg/L	Result µg/L
Acenaphthene	2	U	N-nitrosodimethylamine	2	U
Acenaphthylene	2	U	N-nitroso-di-n-propylamine	2	U
Anthracene	2	U	n-nitrosodiphenylamine	2	U
Benzo[a]anthracene	2	U	Pyridine	2	U
Benzo[a] pyrene	2	U	2-Methylnaphthalene	2	U
Benzo[b] fluoranthene	2	U	2-Chloronaphthalene	2	U
Benzo[k] fluoranthene	2	U	Naphthalene	2	U
Benzo(g,h,i) perylene	2	U	Phenanthrene	2	U
Chrysene	2	U	Dibenzofuran	2	U
Dibenz [a,h] anthracene	2	U	Aniline	2	U
Fluoranthene	2	U	4-Chloroaniline	2	U
Fluorene	2	U	2-Nitroaniline	2	U
Indeno [1,2,3-cd] pyrene	2	U	3-Nitroaniline	2	U
Pyrene	2	U	4-Nitroaniline	2	U
Hexachloroethane	2	U	Carbazole	2	U
Isophorone	2	U			

Base Neutral Surrogate Standard Recovery

2-Fluorobiphenyl	83	%	d5-nitrobenzene	81	%	d14-p-terphenyl	73 %

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis was conducted according to: Test Methods for Evaluating Solid Waste, SW-846 Method 8270C.

COMMENTS: *Due to the reactive nature of this compound, the Benzidine quantitation limit is estimated.

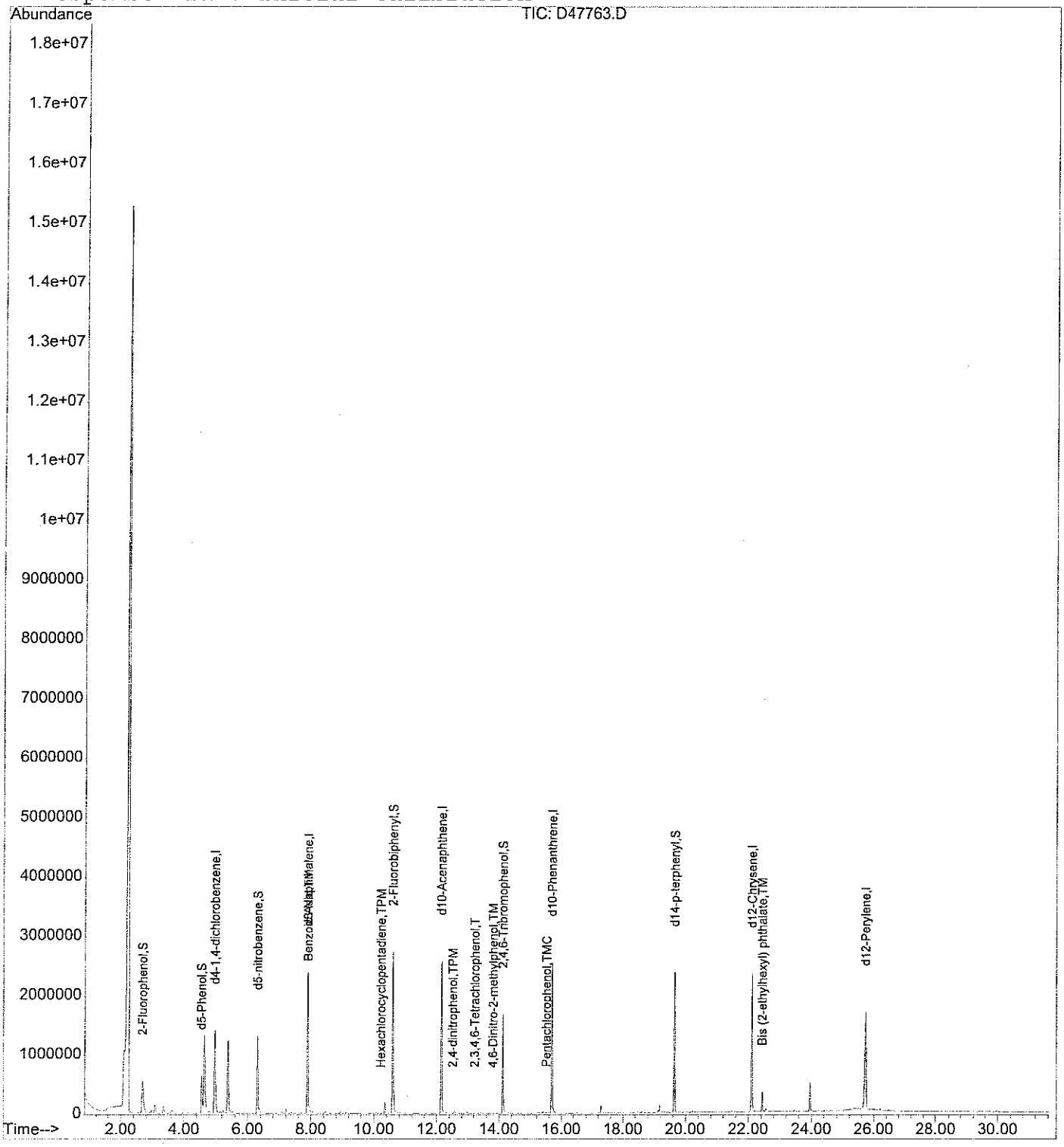
8270/525 layout

Authorized signature

Quantitation Report

Data File : D:\HPCHEM\1\DATA\053006-D\D47763.D Vial: 9
Acq On : 30 May 2006 12:35 pm Operator:
Sample : 56637-3 Inst : D Instrum
Misc : Multiplr: 1.00
MS Integration Params: RTEINT.P
Quant Time: May 30 12:38 2006 Quant Results File: S05246.RES

Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Tue May 30 09:52:41 2006
Response via : Initial Calibration



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May 31, 2006

CLIENT SAMPLE ID

Project Name: GP Old Town Mill

Project Number: 218736.002

Field Sample ID: SB-22 (0-4)

SAMPLE DATA

Lab Sample ID: 56637-4
Matrix: Solid
Percent Solid: 85
Dilution Factor: 1.2
Collection Date: 05/25/06
Lab Receipt Date: 05/27/06
Extraction Date: 05/30/06
Analysis Date: 05/31/06

ANALYTICAL RESULTS SEMI-VOLATILE ORGANICS

ACID COMPOUND	Quantitation Limit µg/kg	Result µg/kg	ACID COMPOUND	Quantitation Limit µg/kg	Result µg/kg
2-Chlorophenol	290	U	Pentachlorophenol	290	U
4-Chloro-3-methylphenol	290	U	Phenol	290	U
2,4-Dichlorophenol	290	U	2,4,5-Trichlorophenol	290	U
2,4-Dimethylphenol	290	U	2,4,6-Trichlorophenol	290	U
2,4-dinitrophenol	290	U	Benzoic Acid	290	U
4,6-Dinitro-2-methylphenol	290	U	2-Methylphenol	290	U
2-Nitrophenol	290	U	3+4-Methylphenol	290	U
2,6-Dichlorophenol	290	U	Benzyl Alcohol	290	U
4-Nitrophenol	290	U	2,3,4,6-Tetrachlorophenol	290	U

Acid Surrogate Standard Recovery

2-Fluorophenol	68 %	d5-Phenol	80 %	2,4,6-Tribromophenol	90 %

BASE NEUTRAL COMPOUND	Quantitation Limit µg/kg	Result µg/kg	BASE NEUTRAL COMPOUND	Quantitation Limit µg/kg	Result µg/kg
1,2-Dichlorobenzene	290	U	Hexachlorobenzene	290	U
1,3-Dichlorobenzene	290	U	* Benzidine	290	U
1,4-Dichlorobenzene	290	U	3,3'-Dichlorobenzidine	290	U
2,4-Dinitrotoluene	290	U	Azobenzene	290	U
2,6-Dinitrotoluene	290	U	Bis(2-chloroethoxy)methane	290	U
Nitrobenzene	290	U	bis(2-chloroethyl) ether	290	U
Hexachlorobutadiene	290	U	bis(2-chloroisopropyl)ether	290	U
Dimethyl Phthalate	290	U	4-bromophenyl phenyl ether	290	U
Di-n-butyl phthalate	290	U	Butyl benzyl phthalate	290	U
di-n-octyl-phthalate	290	U	4-Chlorophenyl phenyl ether	290	U
Bis (2-ethylhexyl) phthalate	290	U	Diethyl Phthalate	290	U
1,2,4-Trichlorobenzene	290	U	Hexachlorocyclopentadiene	290	U

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis was conducted according to: Test Methods for Evaluating Solid Waste, SW-846 Method 8270C.

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8270/625 layout

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Portland ME 04102

May 31, 2006

CLIENT SAMPLE ID

Project Name: GP Old Town Mill

Project Number: 218736.002

Field Sample ID: SB-22 (0-4)

SAMPLE DATA

Lab Sample ID: 56637-4
Matrix: Solid
Percent Solid: 85
Dilution Factor: 1.2
Collection Date: 05/25/06
Lab Receipt Date: 05/27/06
Extraction Date: 05/30/06
Analysis Date: 05/31/06

ANALYTICAL RESULTS SEMI-VOLATILE ORGANICS

BASE NEUTRAL COMPOUND	Quantitation Limit µg/kg	Result µg/kg	BASE NEUTRAL COMPOUND	Quantitation Limit µg/kg	Result µg/kg
Acenaphthene	290	U	N-nitrosodimethylamine	290	U
Acenaphthylene	290	U	N-nitroso-di-n-propylamine	290	U
Anthracene	290	U	n-nitrosodiphenylamine	290	U
Benzo[a]anthracene	290	U	Pyridine	290	U
Benzo[a] pyrene	290	U	2-Methylnaphthalene	290	U
Benzo[b] fluoranthene	290	U	2-Chloronaphthalene	290	U
Benzo[k] fluoranthene	290	U	Naphthalene	290	U
Benzo(g,h,i) perylene	290	U	Phenanthrene	290	U
Chrysene	290	U	Dibenzofuran	290	U
Dibenz [a,h] anthracene	290	U	Aniline	290	U
Fluoranthene	290	U	4-Chloroaniline	290	U
Fluorene	290	U	2-Nitroaniline	290	U
Indeno [1,2,3-cd] pyrene	290	U	3-Nitroaniline	290	U
Pyrene	290	U	4-Nitroaniline	290	U
Hexachloroethane	290	U	Carbazole	290	U
Isophorone	290	U			

Base Neutral Surrogate Standard Recovery

2-Fluorobiphenyl	77	%	d5-nitrobenzene	74	%	d14-p-terphenyl	93 %

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis was conducted according to: Test Methods for Evaluating Solid Waste, SW-846 Method 8270C.

COMMENTS: *Due to the reactive nature of this compound, the Benzidine quantitation limit is estimated.
Results are expressed on a dry weight basis.

8270/625 layout

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May 31, 2006

CLIENT SAMPLE ID

Project Name: GP Old Town Mill

Project Number: 218736.002

Field Sample ID: SB-21 (0-4')

SAMPLE DATA

Lab Sample ID: 56637-9
Matrix: Solid
Percent Solid: 97
Dilution Factor: 1.0
Collection Date: 05/25/06
Lab Receipt Date: 05/27/06
Extraction Date: 05/30/06
Analysis Date: 05/31/06

ANALYTICAL RESULTS SEMI-VOLATILE ORGANICS

ACID COMPOUND	Quantitation Limit µg/kg	Result µg/kg	ACID COMPOUND	Quantitation Limit µg/kg	Result µg/kg
2-Chlorophenol	250	U	Pentachlorophenol	250	U
4-Chloro-3-methylphenol	250	U	Phenol	250	U
2,4-Dichlorophenol	250	U	2,4,5-Trichlorophenol	250	U
2,4-Dimethylphenol	250	U	2,4,6-Trichlorophenol	250	U
2,4-dinitrophenol	250	U	Benzoic Acid	250	U
4,6-Dinitro-2-methylphenol	250	U	2-Methylphenol	250	U
2-Nitrophenol	250	U	3+4-Methylphenol	250	U
2,6-Dichlorophenol	250	U	Benzyl Alcohol	250	U
4-Nitrophenol	250	U	2,3,4,6-Tetrachlorophenol	250	U

Acid Surrogate Standard Recovery

2-Fluorophenol	61 %	d5-Phenol	72 %	2,4,6-Tribromophenol	83 %

BASE NEUTRAL COMPOUND	Quantitation Limit µg/kg	Result µg/kg	BASE NEUTRAL COMPOUND	Quantitation Limit µg/kg	Result µg/kg
1,2-Dichlorobenzene	250	U	Hexachlorobenzene	250	U
1,3-Dichlorobenzene	250	U	* Benzidine	250	U
1,4-Dichlorobenzene	250	U	3,3'-Dichlorobenzidine	250	U
2,4-Dinitrotoluene	250	U	Azobenzene	250	U
2,6-Dinitrotoluene	250	U	Bis(2-chloroethoxy)methane	250	U
Nitrobenzene	250	U	bis(2-chloroethyl) ether	250	U
Hexachlorobutadiene	250	U	bis(2-chloroisopropyl)ether	250	U
Dimethyl Phthalate	250	U	4-bromophenyl phenyl ether	250	U
Di-n-butyl phthalate	250	U	Butyl benzyl phthalate	250	U
di-n-octyl-phthalate	250	U	4-Chlorophenyl phenyl ether	250	U
Bis (2-ethylhexyl) phthalate	250	U	Diethyl Phthalate	250	U
1,2,4-Trichlorobenzene	250	U	Hexachlorocyclopentadiene	250	U

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis was conducted according to: Test Methods for Evaluating Solid Waste, SW-846 Method 8270C.

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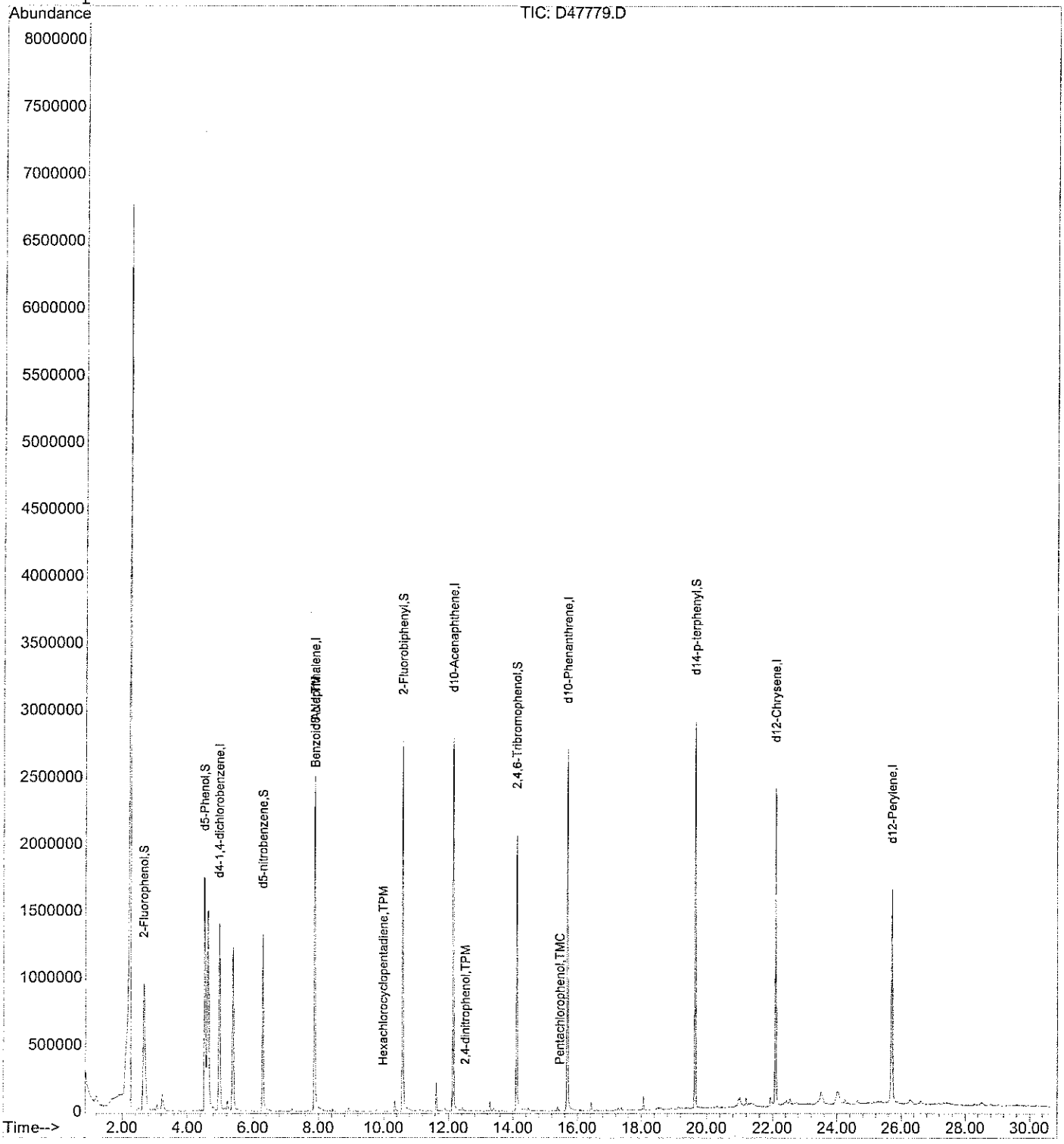
8270/625 layout

Quantitation Report

Data File : D:\HPCHEM\1\DATA\053006-D\D47779.D Vial: 9
Acq On : 31 May 2006 1:23 am Operator:
Sample : 56637-4 Inst : D Instrum
Misc : SOIL Multiplr: 1.00
MS Integration Params: RTEINT.P
Quant Time: May 31 8:21 2006 Quant Results File: S05246.RES

Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Tue May 30 20:31:56 2006
Response via : Initial Calibration



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May 31, 2006

CLIENT SAMPLE ID

Project Name: GP Old Town Mill

Project Number: 218736.002

Field Sample ID: SB-21 (0-4')

SAMPLE DATA

Lab Sample ID: 56637-9
Matrix: Solid
Percent Solid: 97
Dilution Factor: 1.0
Collection Date: 05/25/06
Lab Receipt Date: 05/27/06
Extraction Date: 05/30/06
Analysis Date: 05/31/06

ANALYTICAL RESULTS SEMI-VOLATILE ORGANICS

BASE NEUTRAL COMPOUND	Quantitation Limit µg/kg	Result µg/kg	BASE NEUTRAL COMPOUND	Quantitation Limit µg/kg	Result µg/kg
Acenaphthene	250	U	N-nitrosodimethylamine	250	U
Acenaphthylene	250	U	N-nitroso-di-n-propylamine	250	U
Anthracene	250	U	n-nitrosodiphenylamine	250	U
Benzo[a]anthracene	250	U	Pyridine	250	U
Benzo[a] pyrene	250	U	2-Methylnaphthalene	250	U
Benzo[b] fluoranthene	250	U	2-Chloronaphthalene	250	U
Benzo[k] fluoranthene	250	U	Naphthalene	250	U
Benzo(g,h,i) perylene	250	U	Phenanthrene	250	U
Chrysene	250	U	Dibenzofuran	250	U
Dibenz [a,h] anthracene	250	U	Aniline	250	U
Fluoranthene	250	U	4-Chloroaniline	250	U
Fluorene	250	U	2-Nitroaniline	250	U
Indeno [1,2,3-cd] pyrene	250	U	3-Nitroaniline	250	U
Pyrene	250	U	4-Nitroaniline	250	U
Hexachloroethane	250	U	Carbazole	250	U
Isophorone	250	U			

Base Neutral Surrogate Standard Recovery

2-Fluorobiphenyl	72	%	d5-nitrobenzene	66	%	d14-p-terphenyl	85	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis was conducted according to: Test Methods for Evaluating Solid Waste, SW-846 Method 8270C.

COMMENTS: *Due to the reactive nature of this compound, the Benzidine quantitation limit is estimated.
Results are expressed on a dry weight basis.

8270/625 layout

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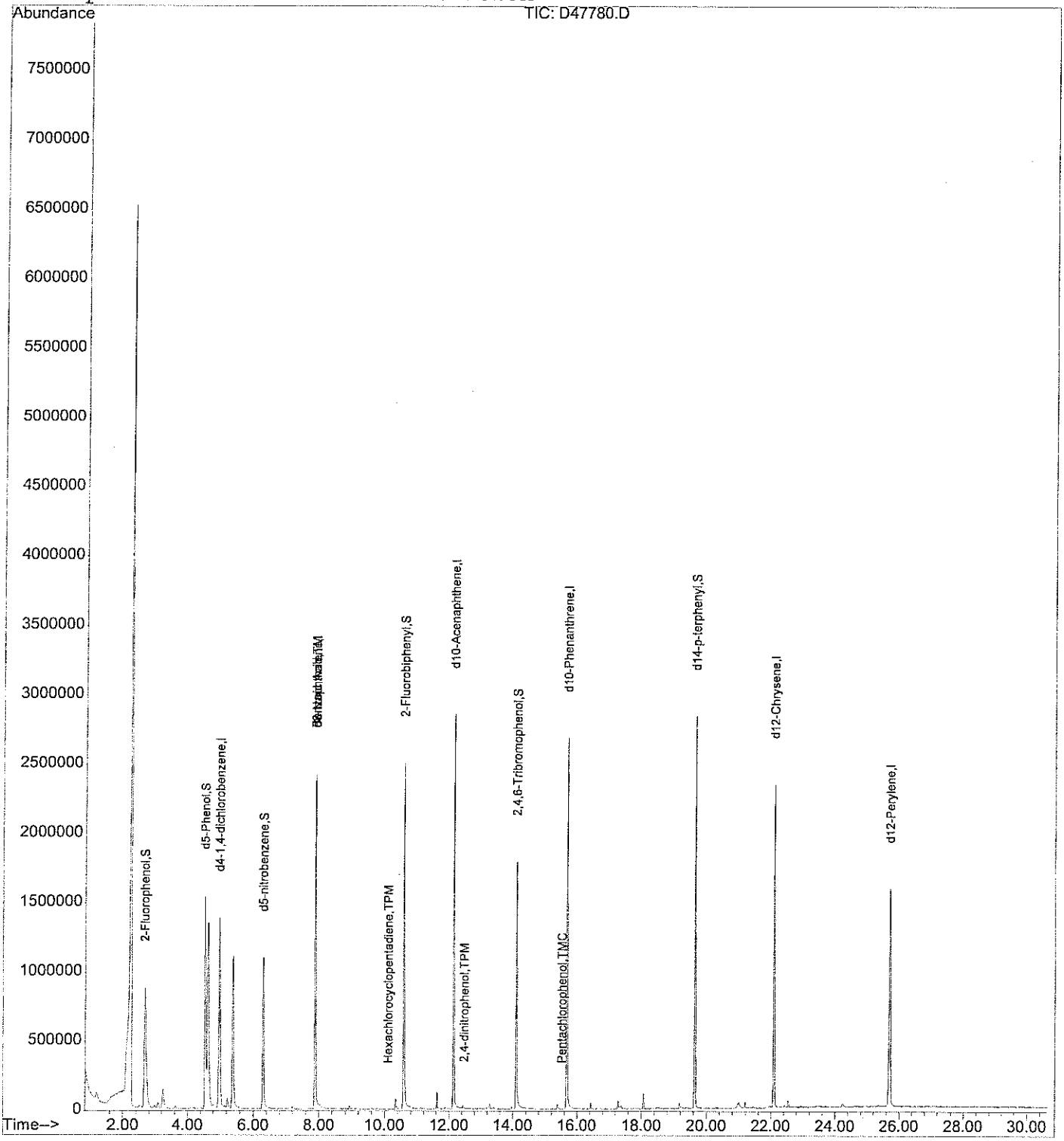
Quantitation Report

Data File : D:\HPCHEM\1\DATA\053006-D\D47780.D Vial: 10
Acq On : 31 May 2006 1:58 am Operator:
Sample : 56637-9 Inst : D Instrum
Misc : SOIL Multiplr: 1.00
MS Integration Params: RTEINT.P
Quant Time: May 31 8:21 2006 Quant Results File: S05246.RES

Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Tue May 30 20:31:56 2006
Response via : Initial Calibration



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May 31, 2006

CLIENT SAMPLE ID

Project Name: GP Old Town Mill

Project Number: 218736.002

Field Sample ID: SB-5 (0-4')

SAMPLE DATA

Lab Sample ID: 56637-10
Matrix: Solid
Percent Solid: 94
Dilution Factor: 1.0
Collection Date: 05/25/06
Lab Receipt Date: 05/27/06
Extraction Date: 05/30/06
Analysis Date: 05/31/06

ANALYTICAL RESULTS SEMI-VOLATILE ORGANICS

ACID COMPOUND	Quantitation Limit µg/kg	Result µg/kg	ACID COMPOUND	Quantitation Limit µg/kg	Result µg/kg
2-Chlorophenol	260	U	Pentachlorophenol	260	U
4-Chloro-3-methylphenol	260	U	Phenol	260	U
2,4-Dichlorophenol	260	U	2,4,5-Trichlorophenol	260	U
2,4-Dimethylphenol	260	U	2,4,6-Trichlorophenol	260	U
2,4-dinitrophenol	260	U	Benzoic Acid	260	U
4,6-Dinitro-2-methylphenol	260	U	2-Methylphenol	260	U
2-Nitrophenol	260	U	3+4-Methylphenol	260	U
2,6-Dichlorophenol	260	U	Benzyl Alcohol	260	U
4-Nitrophenol	260	U	2,3,4,6-Tetrachlorophenol	260	U

Acid Surrogate Standard Recovery

2-Fluorophenol	56 %	d5-Phenol	67 %	2,4,6-Tribromophenol	87 %

BASE NEUTRAL COMPOUND	Quantitation Limit µg/kg	Result µg/kg	BASE NEUTRAL COMPOUND	Quantitation Limit µg/kg	Result µg/kg
1,2-Dichlorobenzene	260	U	Hexachlorobenzene	260	U
1,3-Dichlorobenzene	260	U	* Benzidine	260	U
1,4-Dichlorobenzene	260	U	3,3'-Dichlorobenzidine	260	U
2,4-Dinitrotoluene	260	U	Azobenzene	260	U
2,6-Dinitrotoluene	260	U	Bis(2-chloroethoxy)methane	260	U
Nitrobenzene	260	U	bis(2-chloroethyl) ether	260	U
Hexachlorobutadiene	260	U	bis(2-chloroisopropyl)ether	260	U
Dimethyl Phthalate	260	U	4-bromophenyl phenyl ether	260	U
Di-n-butyl phthalate	260	U	Butyl benzyl phthalate	260	U
di-n-octyl-phthalate	260	U	4-Chlorophenyl phenyl ether	260	U
Bis (2-ethylhexyl) phthalate	260	U	Diethyl Phthalate	260	U
1,2,4-Trichlorobenzene	260	U	Hexachlorocyclopentadiene	260	U

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis was conducted according to: Test Methods for Evaluating Solid Waste, SW-846 Method 8270C.

Authorized signature Milena Tulli

8270/625 layout

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May 31, 2006

CLIENT SAMPLE ID

Project Name: GP Old Town Mill

Project Number: 218736.002

Field Sample ID: SB-5 (0-4')

SAMPLE DATA

Lab Sample ID: 56637-10
Matrix: Solid
Percent Solid: 94
Dilution Factor: 1.0
Collection Date: 05/25/06
Lab Receipt Date: 05/27/06
Extraction Date: 05/30/06
Analysis Date: 05/31/06

ANALYTICAL RESULTS SEMI-VOLATILE ORGANICS

BASE NEUTRAL COMPOUND	Quantitation Limit µg/kg	Result µg/kg	BASE NEUTRAL COMPOUND	Quantitation Limit µg/kg	Result µg/kg
Acenaphthene	260	U	N-nitrosodimethylamine	260	U
Acenaphthylene	260	U	N-nitroso-di-n-propylamine	260	U
Anthracene	260	U	n-nitrosodiphenylamine	260	U
Benzo[a]anthracene	260	U	Pyridine	260	U
Benzo[a] pyrene	260	U	2-Methylnaphthalene	260	U
Benzo[b] fluoranthene	260	U	2-Chloronaphthalene	260	U
Benzo[k] fluoranthene	260	U	Naphthalene	260	U
Benzo(g,h,i) perylene	260	U	Phenanthrene	260	U
Chrysene	260	U	Dibenzofuran	260	U
Dibenz [a,h] anthracene	260	U	Aniline	260	U
Fluoranthene	260	U	4-Chloroaniline	260	U
Fluorene	260	U	2-Nitroaniline	260	U
Indeno [1,2,3-cd] pyrene	260	U	3-Nitroaniline	260	U
Pyrene	260	U	4-Nitroaniline	260	U
Hexachloroethane	260	U	Carbazole	260	U
Isophorone	260	U			

Base Neutral Surrogate Standard Recovery

2-Fluorobiphenyl	68	%	d5-nitrobenzene	59	%	d14-p-terphenyl	88 %

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis was conducted according to: Test Methods for Evaluating Solid Waste, SW-846 Method 8270C.

COMMENTS: *Due to the reactive nature of this compound, the Benzidine quantitation limit is estimated.
Results are expressed on a dry weight basis.

8270/625 layout

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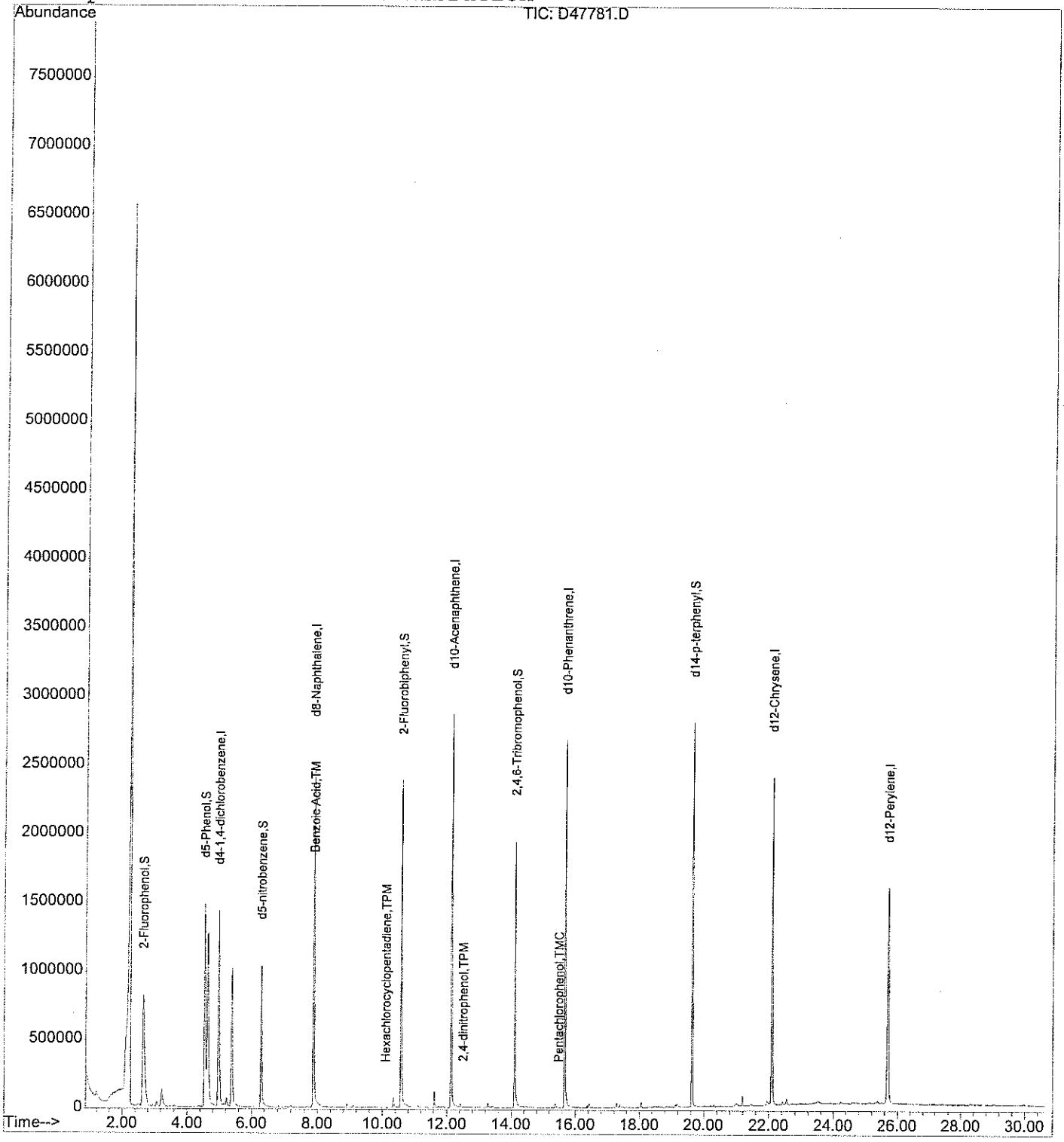
Quantitation Report

Data File : D:\HPCHEM\1\DATA\053006-D\D47781.D Vial: 11
Acq On : 31 May 2006 2:31 am Operator:
Sample : 56637-10 Inst : D Instrum
Misc : SOIL Multiplr: 1.00
MS Integration Params: RTEINT.P
Quant Time: May 31 8:21 2006 Quant Results File: S05246.RES

Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Tue May 30 20:31:56 2006
Response via : Initial Calibration



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May 31, 2006

CLIENT SAMPLE ID

Project Name: GP Old Town Mill

Project Number: 218736.002

Field Sample ID: SB-23 (0-4')

SAMPLE DATA

Lab Sample ID: 56637-12
Matrix: Solid
Percent Solid: 90
Dilution Factor: 1.1
Collection Date: 05/25/06
Lab Receipt Date: 05/27/06
Extraction Date: 05/30/06
Analysis Date: 05/31/06

ANALYTICAL RESULTS SEMI-VOLATILE ORGANICS

ACID COMPOUND	Quantitation Limit µg/kg	Result µg/kg	ACID COMPOUND	Quantitation Limit µg/kg	Result µg/kg
2-Chlorophenol	280	U	Pentachlorophenol	280	U
4-Chloro-3-methylphenol	280	U	Phenol	280	U
2,4-Dichlorophenol	280	U	2,4,5-Trichlorophenol	280	U
2,4-Dimethylphenol	280	U	2,4,6-Trichlorophenol	280	U
2,4-dinitrophenol	280	U	Benzoic Acid	280	U
4,6-Dinitro-2-methylphenol	280	U	2-Methylphenol	280	U
2-Nitrophenol	280	U	3+4-Methylphenol	280	U
2,6-Dichlorophenol	280	U	Benzyl Alcohol	280	U
4-Nitrophenol	280	U	2,3,4,6-Tetrachlorophenol	280	U

Acid Surrogate Standard Recovery

2-Fluorophenol	67 %	d5-Phenol	77 %	2,4,6-Tribromophenol	80 %

BASE NEUTRAL COMPOUND	Quantitation Limit µg/kg	Result µg/kg	BASE NEUTRAL COMPOUND	Quantitation Limit µg/kg	Result µg/kg
1,2-Dichlorobenzene	280	U	Hexachlorobenzene	280	U
1,3-Dichlorobenzene	280	U	* Benzidine	280	U
1,4-Dichlorobenzene	280	U	3,3'-Dichlorobenzidine	280	U
2,4-Dinitrotoluene	280	U	Azobenzene	280	U
2,6-Dinitrotoluene	280	U	Bis(2-chloroethoxy)methane	280	U
Nitrobenzene	280	U	bis(2-chloroethyl) ether	280	U
Hexachlorobutadiene	280	U	bis(2-chloroisopropyl)ether	280	U
Dimethyl Phthalate	280	U	4-bromophenyl phenyl ether	280	U
Di-n-butyl phthalate	280	U	Butyl benzyl phthalate	280	U
di-n-octyl-phthalate	280	U	4-Chlorophenyl phenyl ether	280	U
Bis (2-ethylhexyl) phthalate	280	U	Diethyl Phthalate	280	U
1,2,4-Trichlorobenzene	280	U	Hexachlorocyclopentadiene	280	U

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis was conducted according to: Test Methods for Evaluating Solid Waste, SW-846 Method 8270C.

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8270/625 layout

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May 31, 2006

CLIENT SAMPLE ID

Project Name: GP Old Town Mill

Project Number: 218736.002

Field Sample ID: SB-23 (0-4')

SAMPLE DATA

Lab Sample ID: 56637-12
Matrix: Solid
Percent Solid: 90
Dilution Factor: 1.1
Collection Date: 05/25/06
Lab Receipt Date: 05/27/06
Extraction Date: 05/30/06
Analysis Date: 05/31/06

ANALYTICAL RESULTS SEMI-VOLATILE ORGANICS

BASE NEUTRAL COMPOUND	Quantitation Limit µg/kg	Result µg/kg	BASE NEUTRAL COMPOUND	Quantitation Limit µg/kg	Result µg/kg
Acenaphthene	280	U	N-nitrosodimethylamine	280	U
Acenaphthylene	280	U	N-nitroso-di-n-propylamine	280	U
Anthracene	280	U	n-nitrosodiphenylamine	280	U
Benzo[a]anthracene	280	U	Pyridine	280	U
Benzo[a] pyrene	280	U	2-Methylnaphthalene	280	U
Benzo[b] fluoranthene	280	U	2-Chloronaphthalene	280	U
Benzo[k] fluoranthene	280	U	Naphthalene	280	U
Benzo(g,h,i) perylene	280	U	Phenanthrene	280	U
Chrysene	280	U	Dibenzofuran	280	U
Dibenz [a,h] anthracene	280	U	Aniline	280	U
Fluoranthene	280	U	4-Chloroaniline	280	U
Fluorene	280	U	2-Nitroaniline	280	U
Indeno [1,2,3-cd] pyrene	280	U	3-Nitroaniline	280	U
Pyrene	280	U	4-Nitroaniline	280	U
Hexachloroethane	280	U	Carbazole	280	U
Isophorone	280	U			

Base Neutral Surrogate Standard Recovery

2-Fluorobiphenyl	74	%	d5-nitrobenzene	72	%	d14-p-terphenyl	88	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis was conducted according to: Test Methods for Evaluating Solid Waste, SW-846 Method 8270C.

COMMENTS: *Due to the reactive nature of this compound, the Benzidine quantitation limit is estimated.
Results are expressed on a dry weight basis.

8270/625 layout

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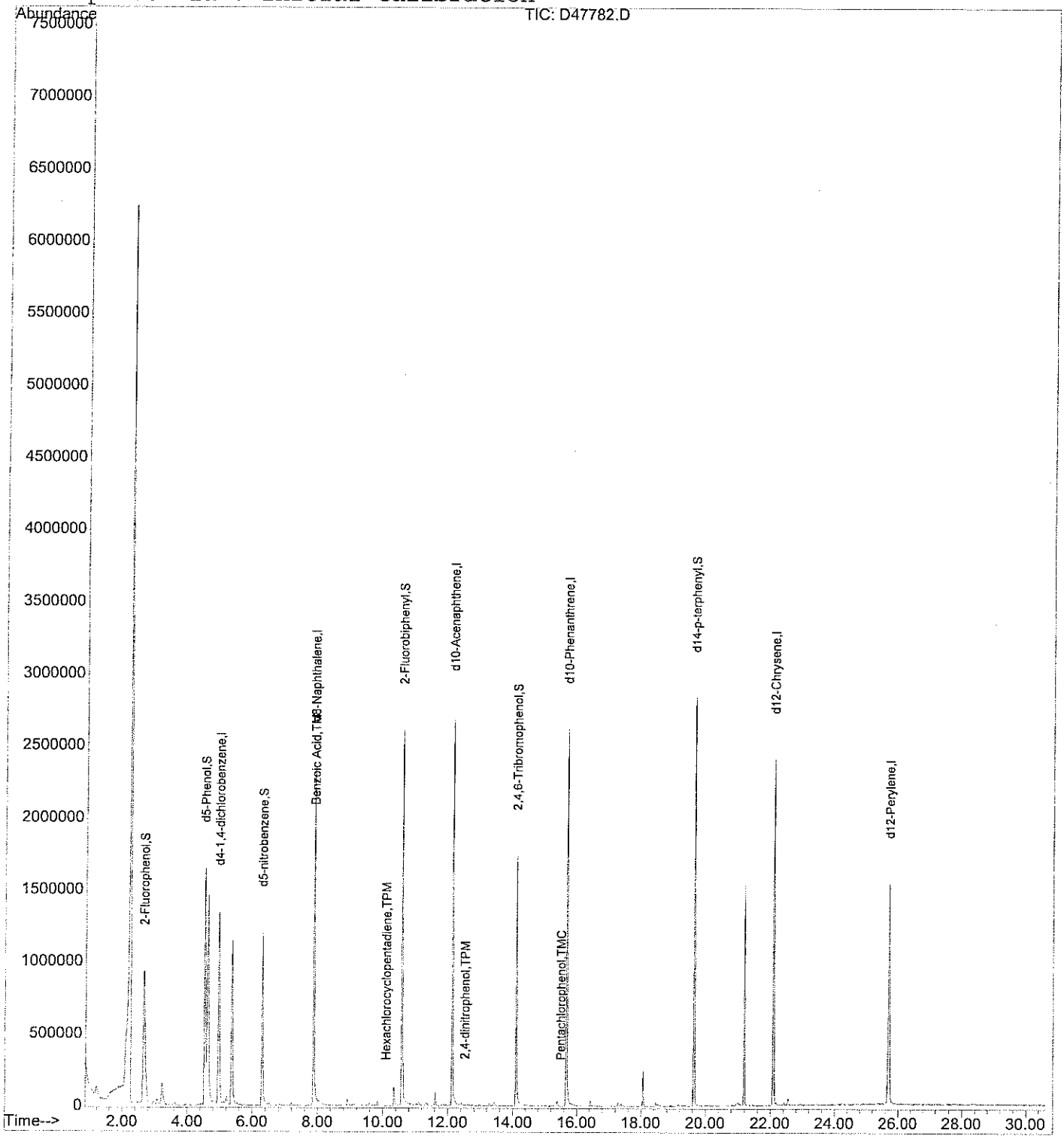
Quantitation Report

Data File : D:\HPCHEM\1\DATA\053006-D\D47782.D Vial: 12
Acq On : 31 May 2006 3:04 am Operator:
Sample : 56637-12 Inst : D Instrum
Misc : SOIL Multiplr: 1.00
MS Integration Params: RTEINT.P
Quant Time: May 31 8:21 2006 Quant Results File: S05246.RES

Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Tue May 30 20:31:56 2006
Response via : Initial Calibration



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May 31, 2006

CLIENT SAMPLE ID

Project Name: GP Old Town Mill

Project Number: 218736.002

Field Sample ID: SB-6 (0-4')

SAMPLE DATA

Lab Sample ID: 56637-11
Matrix: Solid
Percent Solid: 98
Dilution Factor: 1.0
Collection Date: 05/25/06
Lab Receipt Date: 05/27/06
Extraction Date: 05/30/06
Analysis Date: 05/31/06

ANALYTICAL RESULTS SEMI-VOLATILE ORGANICS

ACID COMPOUND	Quantitation Limit µg/kg	Result µg/kg	ACID COMPOUND	Quantitation Limit µg/kg	Result µg/kg
2-Chlorophenol	250	U	Pentachlorophenol	250	U
4-Chloro-3-methylphenol	250	U	Phenol	250	U
2,4-Dichlorophenol	250	U	2,4,5-Trichlorophenol	250	U
2,4-Dimethylphenol	250	U	2,4,6-Trichlorophenol	250	U
2,4-dinitrophenol	250	U	Benzoic Acid	250	U
4,6-Dinitro-2-methylphenol	250	U	2-Methylphenol	250	U
2-Nitrophenol	250	U	3+4-Methylphenol	250	U
2,6-Dichlorophenol	250	U	Benzyl Alcohol	250	U
4-Nitrophenol	250	U	2,3,4,6-Tetrachlorophenol	250	U

Acid Surrogate Standard Recovery

2-Fluorophenol	62 %	d5-Phenol	75 %	2,4,6-Tribromophenol	93 %

BASE NEUTRAL COMPOUND	Quantitation Limit µg/kg	Result µg/kg	BASE NEUTRAL COMPOUND	Quantitation Limit µg/kg	Result µg/kg
1,2-Dichlorobenzene	250	U	Hexachlorobenzene	250	U
1,3-Dichlorobenzene	250	U	* Benzidine	250	U
1,4-Dichlorobenzene	250	U	3,3'-Dichlorobenzidine	250	U
2,4-Dinitrotoluene	250	U	Azobenzene	250	U
2,6-Dinitrotoluene	250	U	Bis(2-chloroethoxy)methane	250	U
Nitrobenzene	250	U	bis(2-chloroethyl) ether	250	U
Hexachlorobutadiene	250	U	bis(2-chloroisopropyl)ether	250	U
Dimethyl Phthalate	250	U	4-bromophenyl phenyl ether	250	U
Di-n-butyl phthalate	250	U	Butyl benzyl phthalate	250	U
di-n-octyl-phthalate	250	U	4-Chlorophenyl phenyl ether	250	U
Bis (2-ethylhexyl) phthalate	250	U	Diethyl Phthalate	250	U
1,2,4-Trichlorobenzene	250	U	Hexachlorocyclopentadiene	250	U

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis was conducted according to: Test Methods for Evaluating Solid Waste, SW-846 Method 8270C.

Authorized signature Mylena Tall

8270/625 layout

analytics environmental laboratory LLC

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800-929-9906

Mr. Karl Kasper
Woodard & Curran
41 Hutchins Drive
Portland ME 04102

May 31, 2006

CLIENT SAMPLE ID

Project Name: GP Old Town Mill

Project Number: 218736.002

Field Sample ID: SB-6 (0-4')

SAMPLE DATA

Lab Sample ID: 56637-11
Matrix: Solid
Percent Solid: 98
Dilution Factor: 1.0
Collection Date: 05/25/06
Lab Receipt Date: 05/27/06
Extraction Date: 05/30/06
Analysis Date: 05/31/06

ANALYTICAL RESULTS SEMI-VOLATILE ORGANICS

BASE NEUTRAL COMPOUND	Quantitation Limit µg/kg	Result µg/kg	BASE NEUTRAL COMPOUND	Quantitation Limit µg/kg	Result µg/kg
Acenaphthene	250	U	N-nitrosodimethylamine	250	U
Acenaphthylene	250	U	N-nitroso-di-n-propylamine	250	U
Anthracene	250	U	n-nitrosodiphenylamine	250	U
Benzo[a]anthracene	250	U	Pyridine	250	U
Benzo[a] pyrene	250	U	2-Methylnaphthalene	250	U
Benzo[b] fluoranthene	250	U	2-Chloronaphthalene	250	U
Benzo[k] fluoranthene	250	U	Naphthalene	250	U
Benzo(g,h,i) perylene	250	U	Phenanthrene	250	U
Chrysene	250	U	Dibenzofuran	250	U
Dibenz [a,h] anthracene	250	U	Aniline	250	U
Fluoranthene	250	U	4-Chloroaniline	250	U
Fluorene	250	U	2-Nitroaniline	250	U
Indeno [1,2,3-cd] pyrene	250	U	3-Nitroaniline	250	U
Pyrene	250	U	4-Nitroaniline	250	U
Hexachloroethane	250	U	Carbazole	250	U
Isophorone	250	U			

Base Neutral Surrogate Standard Recovery

2-Fluorobiphenyl	75	%	d5-nitrobenzene	64	%	d14-p-terphenyl	85 %

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis was conducted according to: Test Methods for Evaluating Solid Waste, SW-846 Method 8270C.

COMMENTS: *Due to the reactive nature of this compound, the Benzidine quantitation limit is estimated.
Results are expressed on a dry weight basis.

8270/625 layout

Authorized signature

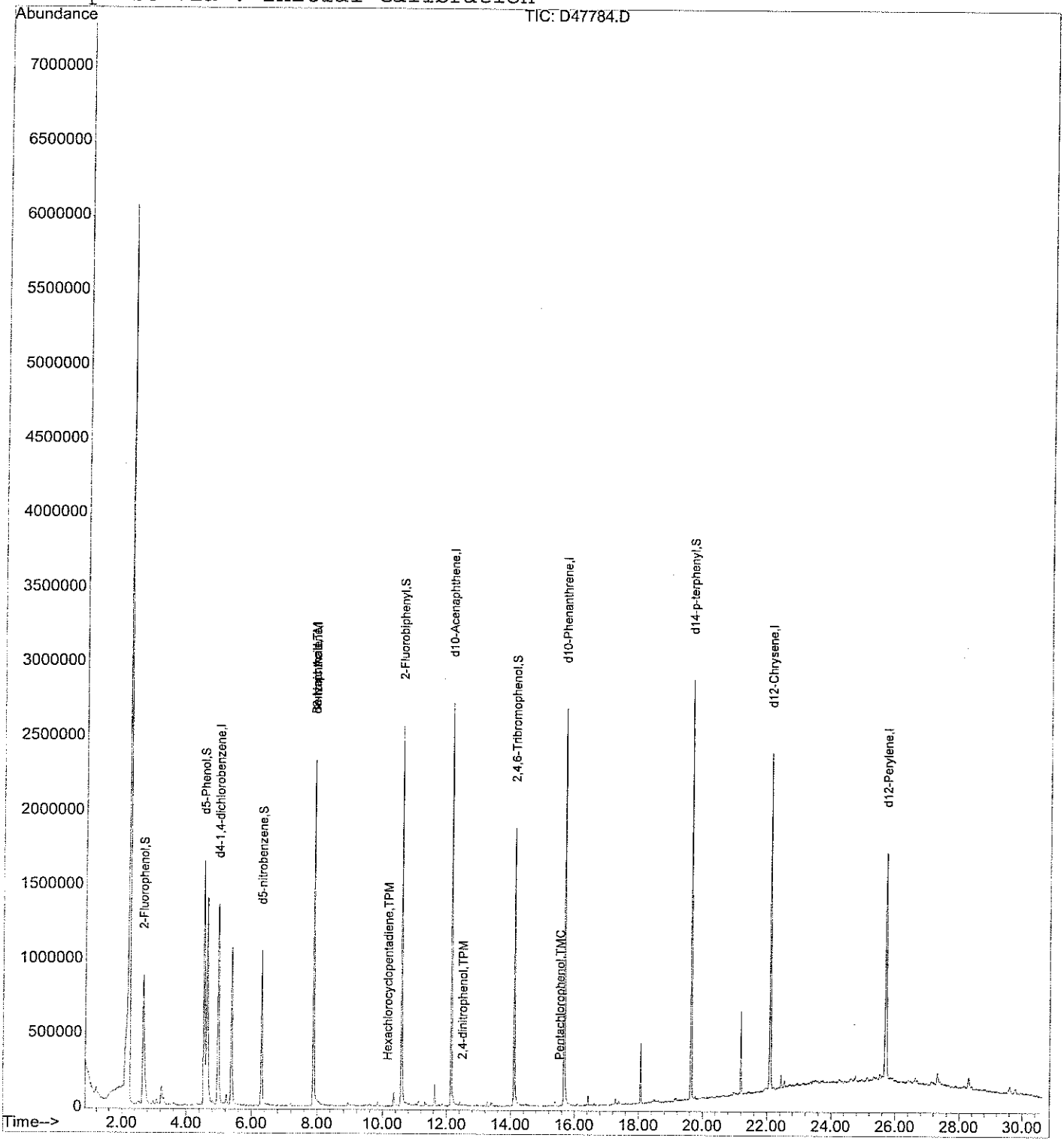
Quantitation Report

Data File : D:\HPCHEM\1\DATA\053006-D\D47784.D
Acq On : 31 May 2006 4:10 am
Sample : 56637-11
Misc : SOIL
MS Integration Params: RTEINT.P
Quant Time: May 31 8:21 2006

Vial: 14
Operator:
Inst : D Instrum
Multiplr: 1.00

Quant Results File: S05246.RES

Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Tue May 30 20:31:56 2006
Response via : Initial Calibration



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Mr. Karl Kasper
Woodard & Curran
41 Hutchins Drive
Portland ME 04102

May 31, 2006

CLIENT SAMPLE ID

Project Name: GP Old Town Mill

Project Number: 218736.002

Field Sample ID: SB-31 (0-4)

SAMPLE DATA

Lab Sample ID: 56637-14
Matrix: Solid
Percent Solid: 63
Dilution Factor: 1.6
Collection Date: 05/25/06
Lab Receipt Date: 05/27/06
Extraction Date: 05/30/06
Analysis Date: 05/31/06

ANALYTICAL RESULTS SEMI-VOLATILE ORGANICS

ACID COMPOUND	Quantitation Limit µg/kg	Result µg/kg	ACID COMPOUND	Quantitation Limit µg/kg	Result µg/kg
2-Chlorophenol	390	U	Pentachlorophenol	390	U
4-Chloro-3-methylphenol	390	U	Phenol	390	U
2,4-Dichlorophenol	390	U	2,4,5-Trichlorophenol	390	U
2,4-Dimethylphenol	390	U	2,4,6-Trichlorophenol	390	U
2,4-dinitrophenol	390	U	Benzoic Acid	390	U
4,6-Dinitro-2-methylphenol	390	U	2-Methylphenol	390	U
2-Nitrophenol	390	U	3+4-Methylphenol	390	U
2,6-Dichlorophenol	390	U	Benzyl Alcohol	390	U
4-Nitrophenol	390	U	2,3,4,6-Tetrachlorophenol	390	U

Acid Surrogate Standard Recovery

2-Fluorophenol	65 %	d5-Phenol	77 %	2,4,6-Tribromophenol	90 %

BASE NEUTRAL COMPOUND	Quantitation Limit µg/kg	Result µg/kg	BASE NEUTRAL COMPOUND	Quantitation Limit µg/kg	Result µg/kg
1,2-Dichlorobenzene	390	U	Hexachlorobenzene	390	U
1,3-Dichlorobenzene	390	U	* Benzidine	390	U
1,4-Dichlorobenzene	390	U	3,3'-Dichlorobenzidine	390	U
2,4-Dinitrotoluene	390	U	Azobenzene	390	U
2,6-Dinitrotoluene	390	U	Bis(2-chloroethoxy)methane	390	U
Nitrobenzene	390	U	bis(2-chloroethyl) ether	390	U
Hexachlorobutadiene	390	U	bis(2-chloroisopropyl)ether	390	U
Dimethyl Phthalate	390	U	4-bromophenyl phenyl ether	390	U
Di-n-butyl phthalate	390	U	Butyl benzyl phthalate	390	U
di-n-octyl-phthalate	390	U	4-Chlorophenyl phenyl ether	390	U
Bis (2-ethylhexyl) phthalate	390	U	Diethyl Phthalate	390	U
1,2,4-Trichlorobenzene	390	U	Hexachlorocyclopentadiene	390	U

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis was conducted according to: Test Methods for Evaluating Solid Waste, SW-846 Method 8270C.

Authorized signature

8270/625 layout

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Mr. Karl Kasper
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Portland ME 04102

May 31, 2006

CLIENT SAMPLE ID

Project Name: GP Old Town Mill

Project Number: 218736.002

Field Sample ID: SB-31 (0-4)

SAMPLE DATA

Lab Sample ID: 56637-14
Matrix: Solid
Percent Solid: 63
Dilution Factor: 1.6
Collection Date: 05/25/06
Lab Receipt Date: 05/27/06
Extraction Date: 05/30/06
Analysis Date: 05/31/06

ANALYTICAL RESULTS SEMI-VOLATILE ORGANICS

BASE NEUTRAL COMPOUND	Quantitation Limit µg/kg	Result µg/kg	BASE NEUTRAL COMPOUND	Quantitation Limit µg/kg	Result µg/kg
Acenaphthene	390	U	N-nitrosodimethylamine	390	U
Acenaphthylene	390	U	N-nitroso-di-n-propylamine	390	U
Anthracene	390	**289 J**	n-nitrosodiphenylamine	390	U
Benzo[a]anthracene	390	U	Pyridine	390	U
Benzo[a] pyrene	390	**657**	2-Methylnaphthalene	390	U
Benzo[b] fluoranthene	390	**1070**	2-Chloronaphthalene	390	U
Benzo[k] fluoranthene	390	**1210**	Naphthalene	390	U
Benzo(g,h,i) perylene	390	U	Phenanthrene	390	**1070**
Chrysene	390	**1080**	Dibenzofuran	390	U
Dibenz [a,h] anthracene	390	U	Aniline	390	U
Fluoranthene	390	**1640**	4-Chloroaniline	390	U
Fluorene	390	U	2-Nitroaniline	390	U
Indeno [1,2,3-cd] pyrene	390	U	3-Nitroaniline	390	U
Pyrene	390	**1380**	4-Nitroaniline	390	U
Hexachloroethane	390	U	Carbazole	390	U
Isophorone	390	U			

Base Neutral Surrogate Standard Recovery

2-Fluorobiphenyl	77	%	d5-nitrobenzene	71	%	d14-p-terphenyl	77 %

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis was conducted according to: Test Methods for Evaluating Solid Waste, SW-846 Method 8270C.

COMMENTS: *Due to the reactive nature of this compound, the Benzidine quantitation limit is estimated.
Results are expressed on a dry weight basis. Sample had low recovery of internal standard d12-Perylene. Sample is being reanlayzed to confirm results.

8270/625 layout

Authorized signature Melissa Hall

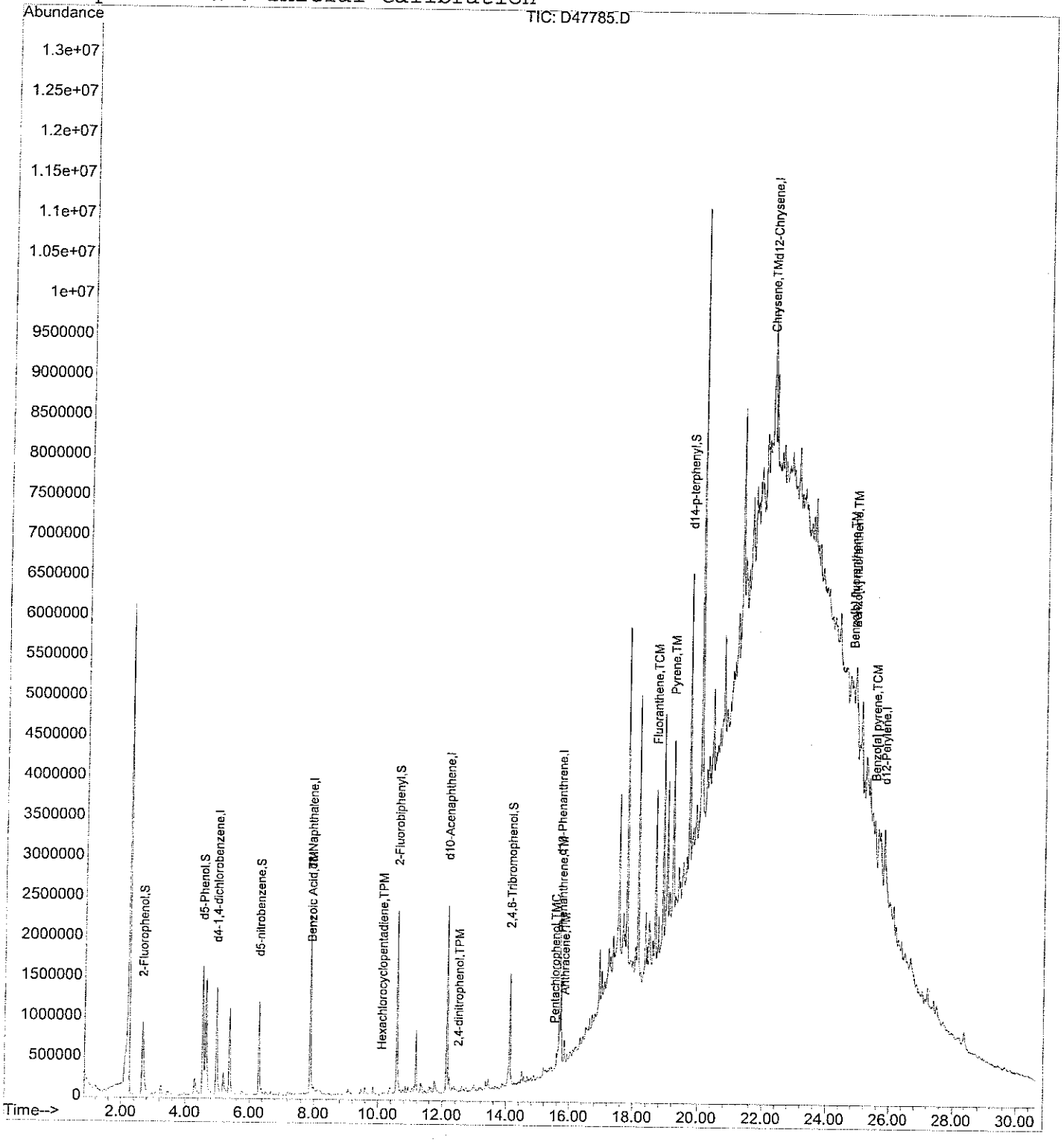
Quantitation Report

Data File : D:\HPCHEM\1\DATA\053006-D\D47785.D
Acq On : 31 May 2006 4:43 am
Sample : 56637-14
Misc : SOIL
MS Integration Params: RTEINT.P
Quant Time: May 31 9:33 2006

Vial: 15
Operator:
Inst : D Instrum
Multiplr: 1.00

Quant Results File: S05246.RES

Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Tue May 30 20:31:56 2006
Response via : Initial Calibration



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Mr. Karl Kasper
Woodard & Curran
41 Hutchins Drive
Portland ME 04102

June 1, 2006

CLIENT SAMPLE ID

Project Name: GP Old Town Mill

Project Number: 218736.002

Field Sample ID: SB-31 (0-4)

SAMPLE DATA

Lab Sample ID: 56637-14 DL
Matrix: Solid
Percent Solid: 63
Dilution Factor: 3.1
Collection Date: 05/25/06
Lab Receipt Date: 05/27/06
Extraction Date: 05/30/06
Analysis Date: 05/31/06

ANALYTICAL RESULTS SEMI-VOLATILE ORGANICS

ACID COMPOUND	Quantitation Limit µg/kg	Result µg/kg	ACID COMPOUND	Quantitation Limit µg/kg	Result µg/kg
2-Chlorophenol	790	U	Pentachlorophenol	790	U
4-Chloro-3-methylphenol	790	U	Phenol	790	U
2,4-Dichlorophenol	790	U	2,4,5-Trichlorophenol	790	U
2,4-Dimethylphenol	790	U	2,4,6-Trichlorophenol	790	U
2,4-dinitrophenol	790	U	Benzoic Acid	790	U
4,6-Dinitro-2-methylphenol	790	U	2-Methylphenol	790	U
2-Nitrophenol	790	U	3+4-Methylphenol	790	U
2,6-Dichlorophenol	790	U	Benzyl Alcohol	790	U
4-Nitrophenol	790	U	2,3,4,6-Tetrachlorophenol	790	U

Acid Surrogate Standard Recovery

2-Fluorophenol	68 %	d5-Phenol	81 %	2,4,6-Tribromophenol	92 %

BASE NEUTRAL COMPOUND	Quantitation Limit µg/kg	Result µg/kg	BASE NEUTRAL COMPOUND	Quantitation Limit µg/kg	Result µg/kg
1,2-Dichlorobenzene	790	U	Hexachlorobenzene	790	U
1,3-Dichlorobenzene	790	U	* Benzidine	790	U
1,4-Dichlorobenzene	790	U	3,3'-Dichlorobenzidine	790	U
2,4-Dinitrotoluene	790	U	Azobenzene	790	U
2,6-Dinitrotoluene	790	U	Bis(2-chloroethoxy)methane	790	U
Nitrobenzene	790	U	bis(2-chloroethyl) ether	790	U
Hexachlorobutadiene	790	U	bis(2-chloroisopropyl)ether	790	U
Dimethyl Phthalate	790	U	4-bromophenyl phenyl ether	790	U
Di-n-butyl phthalate	790	U	Butyl benzyl phthalate	790	U
di-n-octyl-phthalate	790	U	4-Chlorophenyl phenyl ether	790	U
Bis (2-ethylhexyl) phthalate	790	U	Diethyl Phthalate	790	U
1,2,4-Trichlorobenzene	790	U	Hexachlorocyclopentadiene	790	U

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis was conducted according to: Test Methods for Evaluating Solid Waste, SW-846 Method 8270C.

8270/625 layout

Authorized signature

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Mr. Karl Kasper
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41 Hutchins Drive
Portland ME 04102

June 1, 2006

CLIENT SAMPLE ID

Project Name: GP Old Town Mill

Project Number: 218736.002

Field Sample ID: SB-31 (0-4)

SAMPLE DATA

Lab Sample ID:	56637-14	DL
Matrix:	Solid	
Percent Solid:	63	
Dilution Factor:	3.1	
Collection Date:	05/25/06	
Lab Receipt Date:	05/27/06	
Extraction Date:	05/30/06	
Analysis Date:	05/31/06	

ANALYTICAL RESULTS SEMI-VOLATILE ORGANICS

BASE NEUTRAL COMPOUND	Quantitation Limit µg/kg	Result µg/kg	BASE NEUTRAL COMPOUND	Quantitation Limit µg/kg	Result µg/kg
Acenaphthene	790	U	N-nitrosodimethylamine	790	U
Acenaphthylene	790	U	N-nitroso-di-n-propylamine	790	U
Anthracene	790	U	n-nitrosodiphenylamine	790	U
Benzo[a]anthracene	790	**896**	Pyridine	790	U
Benzo[a] pyrene	790	**879**	2-Methylnaphthalene	790	U
Benzo[b] fluoranthene	790	**856**	2-Chloronaphthalene	790	U
Benzo[k] fluoranthene	790	**949**	Naphthalene	790	U
Benzo(g,h,i) perylene	790	U	Phenanthrene	790	**1070**
Chrysene	790	**877**	Dibenzofuran	790	U
Dibenz [a,h] anthracene	790	U	Aniline	790	U
Fluoranthene	790	**1570**	4-Chloroaniline	790	U
Fluorene	790	U	2-Nitroaniline	790	U
Indeno [1,2,3-cd] pyrene	790	U	3-Nitroaniline	790	U
Pyrene	790	**1330**	4-Nitroaniline	790	U
Hexachloroethane	790	U	Carbazole	790	U
Isophorone	790	U			

Base Neutral Surrogate Standard Recovery

2-Fluorobiphenyl	82	%	d5-nitrobenzene	74	%	d14-p-terphenyl	66 %

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis was conducted according to: Test Methods for Evaluating Solid Waste, SW-846 Method 8270C.

COMMENTS: *Due to the reactive nature of this compound, the Benzidine quantitation limit is estimated.
Results are expressed on a dry weight basis.

8270/625 layout

Authorized signature Melissa Tull

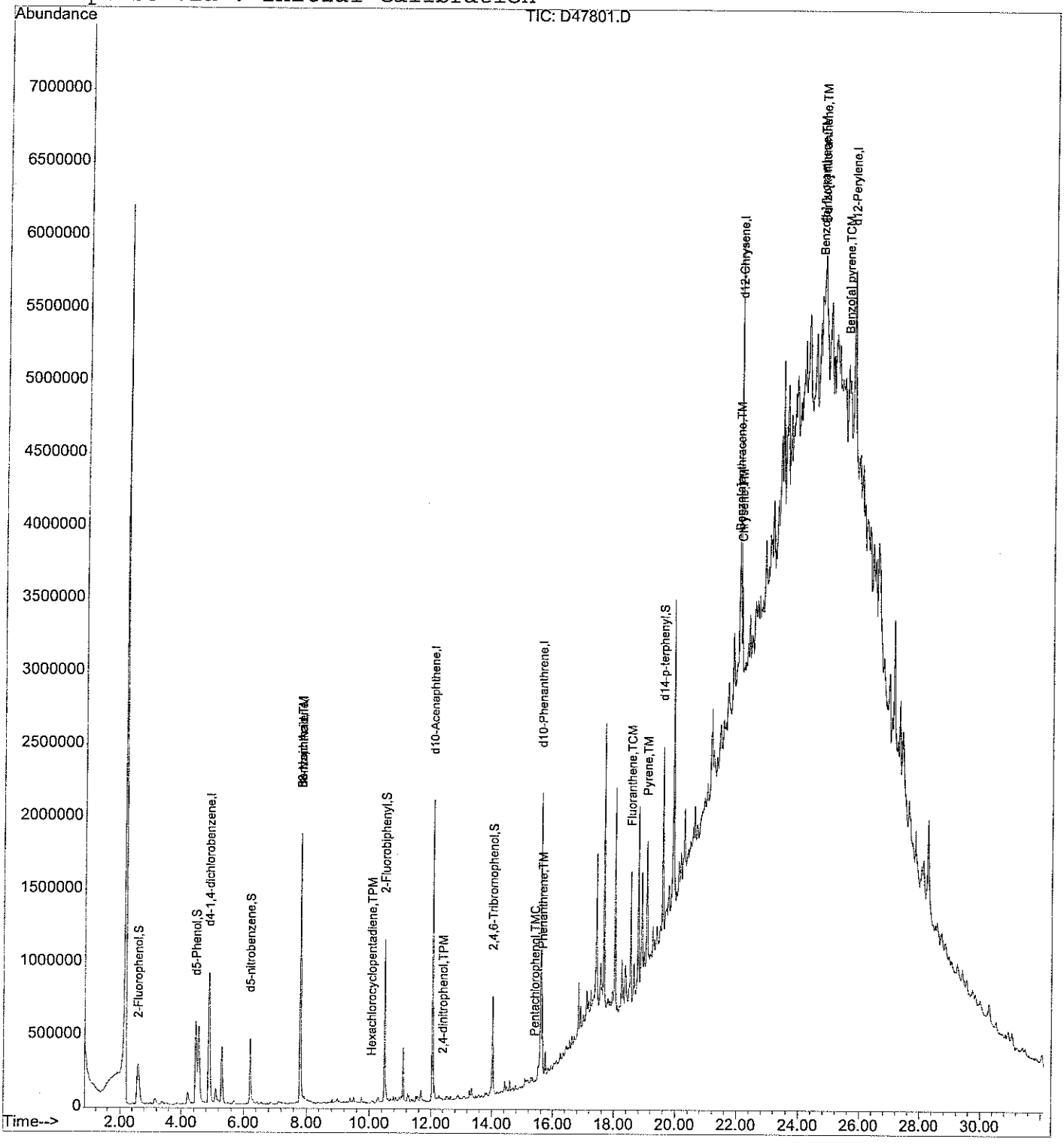
Quantitation Report

Data File : D:\HPCHEM\1\DATA\053106-D\D47801.D
Acq On : 31 May 2006 2:59 pm
Sample : 56637-14,DL,1:2
Misc : SOIL
MS Integration Params: RTEINT.P
Quant Time: Jun 1 1:02 2006

Vial: 8
Operator:
Inst : D Instrum
Multiplr: 1.00

Quant Results File: S05246.RES

Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Wed May 31 19:58:02 2006
Response via : Initial Calibration



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Mr. Karl Kasper
Woodard & Curran
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Portland ME 04102

May 31, 2006

CLIENT SAMPLE ID

Project Name: GP Old Town Mill

Project Number: 218736.002

Field Sample ID: SB-30 (0-4)

SAMPLE DATA

Lab Sample ID: 56637-15
Matrix: Solid
Percent Solid: 81
Dilution Factor: 2.5
Collection Date: 05/25/06
Lab Receipt Date: 05/27/06
Extraction Date: 05/30/06
Analysis Date: 05/31/06

ANALYTICAL RESULTS SEMI-VOLATILE ORGANICS

ACID COMPOUND	Quantitation Limit µg/kg	Result µg/kg	ACID COMPOUND	Quantitation Limit µg/kg	Result µg/kg
2-Chlorophenol	610	U	Pentachlorophenol	610	U
4-Chloro-3-methylphenol	610	U	Phenol	610	U
2,4-Dichlorophenol	610	U	2,4,5-Trichlorophenol	610	U
2,4-Dimethylphenol	610	U	2,4,6-Trichlorophenol	610	U
2,4-dinitrophenol	610	U	Benzoic Acid	610	U
4,6-Dinitro-2-methylphenol	610	U	2-Methylphenol	610	U
2-Nitrophenol	610	U	3+4-Methylphenol	610	U
2,6-Dichlorophenol	610	U	Benzyl Alcohol	610	U
4-Nitrophenol	610	U	2,3,4,6-Tetrachlorophenol	610	U

Acid Surrogate Standard Recovery

2-Fluorophenol	74 %	d5-Phenol	84 %	2,4,6-Tribromophenol	96 %

BASE NEUTRAL COMPOUND	Quantitation Limit µg/kg	Result µg/kg	BASE NEUTRAL COMPOUND	Quantitation Limit µg/kg	Result µg/kg
1,2-Dichlorobenzene	610	U	Hexachlorobenzene	610	U
1,3-Dichlorobenzene	610	U	* Benzidine	610	U
1,4-Dichlorobenzene	610	U	3,3'-Dichlorobenzidine	610	U
2,4-Dinitrotoluene	610	U	Azobenzene	610	U
2,6-Dinitrotoluene	610	U	Bis(2-chloroethoxy)methane	610	U
Nitrobenzene	610	U	bis(2-chloroethyl) ether	610	U
Hexachlorobutadiene	610	U	bis(2-chloroisopropyl)ether	610	U
Dimethyl Phthalate	610	U	4-bromophenyl phenyl ether	610	U
Di-n-butyl phthalate	610	U	Butyl benzyl phthalate	610	U
di-n-octyl-phthalate	610	U	4-Chlorophenyl phenyl ether	610	U
Bis (2-ethylhexyl) phthalate	610	U	Diethyl Phthalate	610	U
1,2,4-Trichlorobenzene	610	U	Hexachlorocyclopentadiene	610	U

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis was conducted according to: Test Methods for Evaluating Solid Waste, SW-846 Method 8270C.

8270/625 layout

Authorized signature Milena Tull

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May 31, 2006

CLIENT SAMPLE ID

Project Name: GP Old Town Mill

Project Number: 218736.002

Field Sample ID: SB-30 (0-4)

SAMPLE DATA

Lab Sample ID: 56637-15
Matrix: Solid
Percent Solid: 81
Dilution Factor: 2.5
Collection Date: 05/25/06
Lab Receipt Date: 05/27/06
Extraction Date: 05/30/06
Analysis Date: 05/31/06

ANALYTICAL RESULTS SEMI-VOLATILE ORGANICS

BASE NEUTRAL COMPOUND	Quantitation Limit µg/kg	Result µg/kg	BASE NEUTRAL COMPOUND	Quantitation Limit µg/kg	Result µg/kg
Acenaphthene	610	**42700 E**	N-nitrosodimethylamine	610	U
Acenaphthylene	610	**318 J**	N-nitroso-di-n-propylamine	610	U
Anthracene	610	**102800 E**	n-nitrosodiphenylamine	610	U
Benzo[a]anthracene	610	U	Pyridine	610	U
Benzo[a] pyrene	610	**34700 E**	2-Methylnaphthalene	610	**2540**
Benzo[b] fluoranthene	610	**92000 E**	2-Chloronaphthalene	610	U
Benzo[k] fluoranthene	610	**49400 E**	Naphthalene	610	**4100**
Benzo(g,h,i) perylene	610	**18500 E**	Phenanthrene	610	U
Chrysene	610	**58700 E**	Dibenzofuran	610	**21900 E**
Dibenz [a,h] anthracene	610	**4360**	Aniline	610	U
Fluoranthene	610	**1230**	4-Chloroaniline	610	U
Fluorene	610	**39300 E**	2-Nitroaniline	610	U
Indeno [1,2,3-cd] pyrene	610	**21500 E**	3-Nitroaniline	610	U
Pyrene	610	**198200 E**	4-Nitroaniline	610	U
Hexachloroethane	610	U	Carbazole	610	U
Isophorone	610	U			

Base Neutral Surrogate Standard Recovery

2-Fluorobiphenyl	79	%	d5-nitrobenzene	75	%	d14-p-terphenyl	220* %

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis was conducted according to: Test Methods for Evaluating Solid Waste, SW-846 Method 8270C.

COMMENTS: *Due to the reactive nature of this compound, the Benzidine quantitation limit is estimated. Results are expressed on a dry weight basis. Sample had low recovery of internal standard d12-Perylene. *Surrogate recovery outside laboratory acceptance criteria. Sample is being reanalyzed at a dilution.

8270/625 layout

Authorized signature

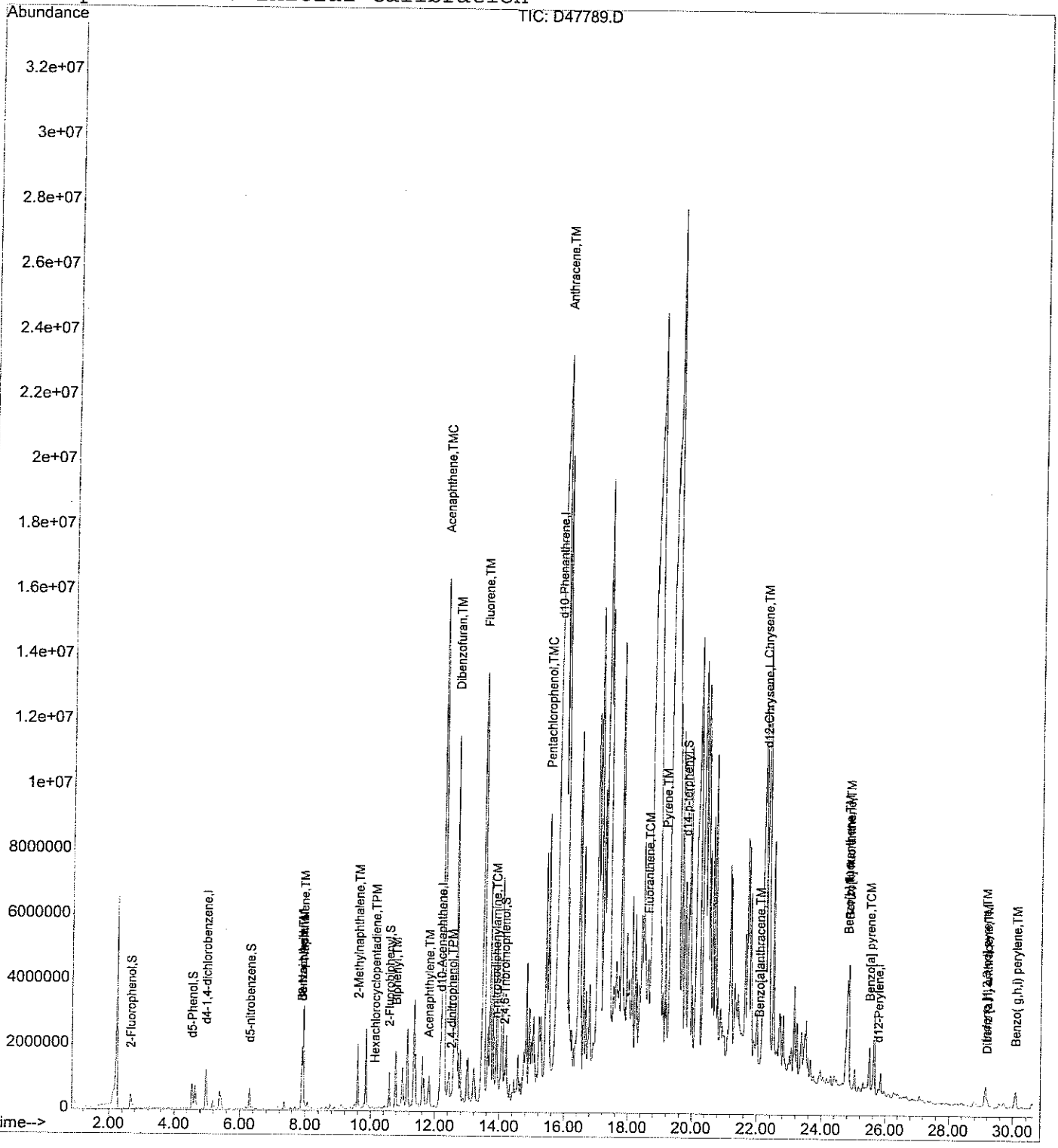
Quantitation Report

Data File : D:\HPCHEM\1\DATA\053006-D\D47789.D
Acq On : 31 May 2006 6:55 am
Sample : 56637-15,,1:2
Misc : SOIL
MS Integration Params: RTEINT.P
Quant Time: May 31 9:47 2006

Vial: 19
Operator:
Inst : D Instrum
Multiplr: 1.00

Quant Results File: S05246.RES

Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Tue May 30 20:31:56 2006
Response via : Initial Calibration



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Mr. Karl Kasper
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Portland ME 04102

June 1, 2006

CLIENT SAMPLE ID

Project Name: GP Old Town Mill

Project Number: 218736.002

Field Sample ID: SB-30 (0-4)

SAMPLE DATA

Lab Sample ID: 56637-15 DL
Matrix: Solid
Percent Solid: 81
Dilution Factor: 123
Collection Date: 05/25/06
Lab Receipt Date: 05/27/06
Extraction Date: 05/30/06
Analysis Date: 05/31/06

ANALYTICAL RESULTS SEMI-VOLATILE ORGANICS

ACID COMPOUND	Quantitation Limit µg/kg	Result µg/kg	ACID COMPOUND	Quantitation Limit µg/kg	Result µg/kg
2-Chlorophenol	31000	U	Pentachlorophenol	31000	U
4-Chloro-3-methylphenol	31000	U	Phenol	31000	U
2,4-Dichlorophenol	31000	U	2,4,5-Trichlorophenol	31000	U
2,4-Dimethylphenol	31000	U	2,4,6-Trichlorophenol	31000	U
2,4-dinitrophenol	31000	U	Benzoic Acid	31000	U
4,6-Dinitro-2-methylphenol	31000	U	2-Methylphenol	31000	U
2-Nitrophenol	31000	U	3+4-Methylphenol	31000	U
2,6-Dichlorophenol	31000	U	Benzyl Alcohol	31000	U
4-Nitrophenol	31000	U	2,3,4,6-Tetrachlorophenol	31000	U

Acid Surrogate Standard Recovery

2-Fluorophenol	* %	d5-Phenol	* %	2,4,6-Tribromophenol	* %

BASE NEUTRAL COMPOUND	Quantitation Limit µg/kg	Result µg/kg	BASE NEUTRAL COMPOUND	Quantitation Limit µg/kg	Result µg/kg
1,2-Dichlorobenzene	31000	U	Hexachlorobenzene	31000	U
1,3-Dichlorobenzene	31000	U	* Benzidine	31000	U
1,4-Dichlorobenzene	31000	U	3,3'-Dichlorobenzidine	31000	U
2,4-Dinitrotoluene	31000	U	Azobenzene	31000	U
2,6-Dinitrotoluene	31000	U	Bis(2-chloroethoxy)methane	31000	U
Nitrobenzene	31000	U	bis(2-chloroethyl) ether	31000	U
Hexachlorobutadiene	31000	U	bis(2-chloroisopropyl)ether	31000	U
Dimethyl Phthalate	31000	U	4-bromophenyl phenyl ether	31000	U
Di-n-butyl phthalate	31000	U	Butyl benzyl phthalate	31000	U
di-n-octyl-phthalate	31000	U	4-Chlorophenyl phenyl ether	31000	U
Bis (2-ethylhexyl) phthalate	31000	U	Diethyl Phthalate	31000	U
1,2,4-Trichlorobenzene	31000	U	Hexachlorocyclopentadiene	31000	U

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis was conducted according to: Test Methods for Evaluating Solid Waste, SW-846 Method 8270C.

8270/625 layout

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Mr. Karl Kasper
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Portland ME 04102

June 1, 2006

CLIENT SAMPLE ID

Project Name: GP Old Town Mill

Project Number: 218736.002

Field Sample ID: SB-30 (0-4)

SAMPLE DATA

Lab Sample ID:	56637-15	DL
Matrix:	Solid	
Percent Solid:	81	
Dilution Factor:	123	
Collection Date:	05/25/06	
Lab Receipt Date:	05/27/06	
Extraction Date:	05/30/06	
Analysis Date:	05/31/06	

ANALYTICAL RESULTS SEMI-VOLATILE ORGANICS

BASE NEUTRAL COMPOUND	Quantitation Limit µg/kg	Result µg/kg	BASE NEUTRAL COMPOUND	Quantitation Limit µg/kg	Result µg/kg
Acenaphthene	31000	**36400**	N-nitrosodimethylamine	31000	U
Acenaphthylene	31000	U	N-nitroso-di-n-propylamine	31000	U
Anthracene	31000	**48900**	n-nitrosodiphenylamine	31000	U
Benzo[a]anthracene	31000	**65300**	Pyridine	31000	U
Benzo[a] pyrene	31000	**52800**	2-Methylnaphthalene	31000	U
Benzo[b] fluoranthene	31000	**46900**	2-Chloronaphthalene	31000	U
Benzo[k] fluoranthene	31000	**44500**	Naphthalene	31000	U
Benzo(g,h,i) perylene	31000	**30500 J**	Phenanthrene	31000	**191300**
Chrysene	31000	**62800**	Dibenzofuran	31000	**17500 J**
Dibenz [a,h] anthracene	31000	U	Aniline	31000	U
Fluoranthene	31000	**187000**	4-Chloroaniline	31000	U
Fluorene	31000	**34700**	2-Nitroaniline	31000	U
Indeno [1,2,3-cd] pyrene	31000	**33700**	3-Nitroaniline	31000	U
Pyrene	31000	**159200**	4-Nitroaniline	31000	U
Hexachloroethane	31000	U	Carbazole	31000	**18500 J**
Isophorone	31000	U			

Base Neutral Surrogate Standard Recovery

2-Fluorobiphenyl	* %	d5-nitrobenzene	* %	d14-p-terphenyl	* %

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis was conducted according to: Test Methods for Evaluating Solid Waste, SW-846 Method 8270C.

COMMENTS: *Due to the reactive nature of this compound, the Benzidine quantitation limit is estimated. Results are expressed on a dry weight basis. *The surrogates were diluted out.

8270/625 layout

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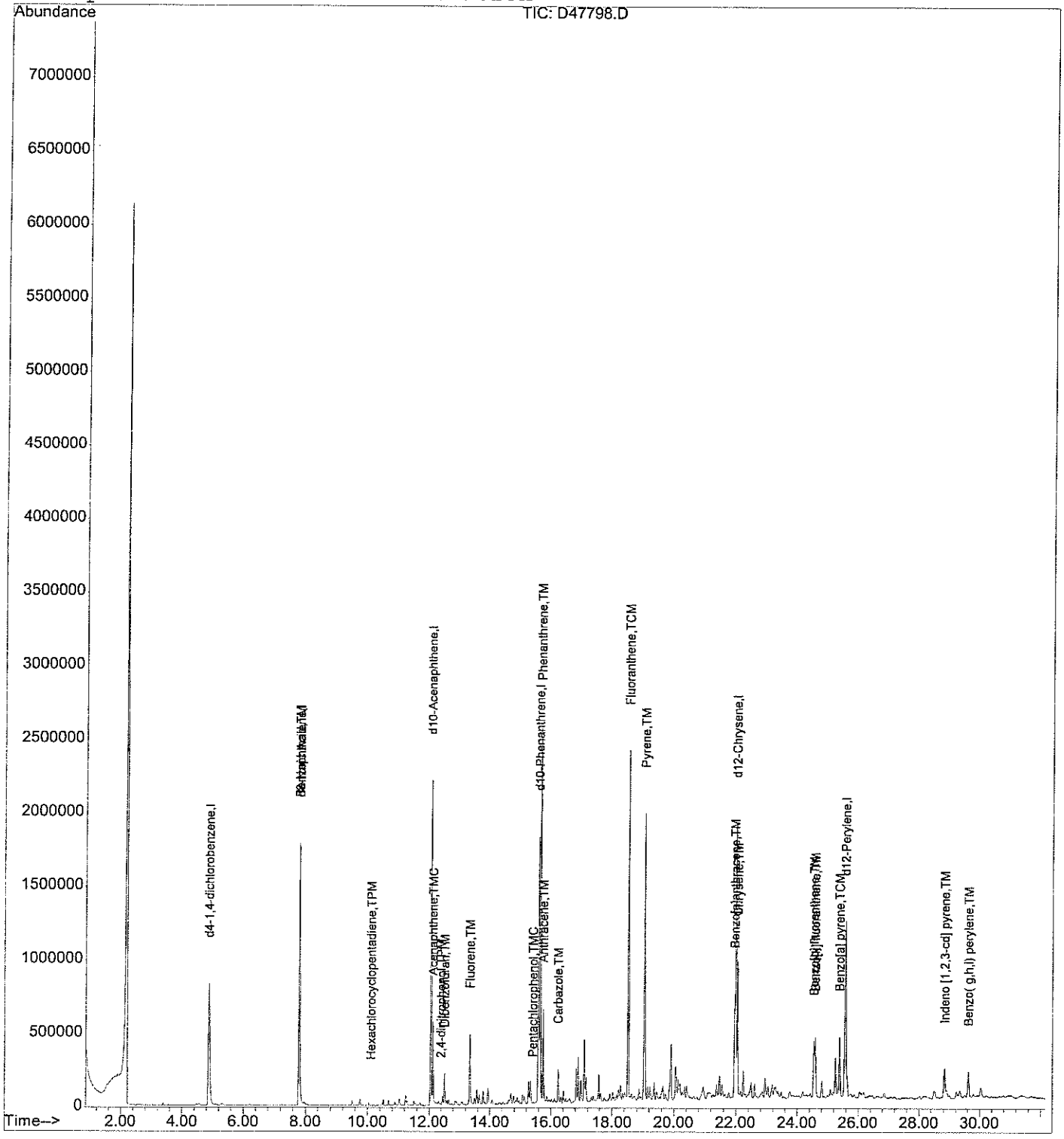
Quantitation Report

Data File : D:\HPCHEM\1\DATA\053106-D\D47798.D
Acq On : 31 May 2006 12:46 pm
Sample : 56637-15,,1:100 DL
Misc : SOIL
MS Integration Params: RTEINT.P
Quant Time: May 31 13:12 2006

Vial: 5
Operator:
Inst : D Instrum
Multiplr: 1.00

Quant Results File: S05246.RES

Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Wed May 31 12:09:53 2006
Response via : Initial Calibration



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Portland ME 04102

May 31, 2006

CLIENT SAMPLE ID

Project Name: GP Old Town Mill

Project Number: 218736.002

Field Sample ID: SB-38 (0-4)

SAMPLE DATA

Lab Sample ID: 56637-16
Matrix: Solid
Percent Solid: 94
Dilution Factor: 5
Collection Date: 05/25/06
Lab Receipt Date: 05/27/06
Extraction Date: 05/30/06
Analysis Date: 05/31/06

ANALYTICAL RESULTS SEMI-VOLATILE ORGANICS

ACID COMPOUND	Quantitation Limit µg/kg	Result µg/kg	ACID COMPOUND	Quantitation Limit µg/kg	Result µg/kg
2-Chlorophenol	1300	U	Pentachlorophenol	1300	U
4-Chloro-3-methylphenol	1300	U	Phenol	1300	U
2,4-Dichlorophenol	1300	U	2,4,5-Trichlorophenol	1300	U
2,4-Dimethylphenol	1300	U	2,4,6-Trichlorophenol	1300	U
2,4-dinitrophenol	1300	U	Benzoic Acid	1300	U
4,6-Dinitro-2-methylphenol	1300	U	2-Methylphenol	1300	U
2-Nitrophenol	1300	U	3+4-Methylphenol	1300	U
2,6-Dichlorophenol	1300	U	Benzyl Alcohol	1300	U
4-Nitrophenol	1300	U	2,3,4,6-Tetrachlorophenol	1300	U

Acid Surrogate Standard Recovery

2-Fluorophenol	66 %	d5-Phenol	74 %	2,4,6-Tribromophenol	79 %

BASE NEUTRAL COMPOUND	Quantitation Limit µg/kg	Result µg/kg	BASE NEUTRAL COMPOUND	Quantitation Limit µg/kg	Result µg/kg
1,2-Dichlorobenzene	1300	U	Hexachlorobenzene	1300	U
1,3-Dichlorobenzene	1300	U	* Benzidine	1300	U
1,4-Dichlorobenzene	1300	U	3,3'-Dichlorobenzidine	1300	U
2,4-Dinitrotoluene	1300	U	Azobenzene	1300	U
2,6-Dinitrotoluene	1300	U	Bis(2-chloroethoxy)methane	1300	U
Nitrobenzene	1300	U	bis(2-chloroethyl) ether	1300	U
Hexachlorobutadiene	1300	U	bis(2-chloroisopropyl)ether	1300	U
Dimethyl Phthalate	1300	U	4-bromophenyl phenyl ether	1300	U
Di-n-butyl phthalate	1300	U	Butyl benzyl phthalate	1300	U
di-n-octyl-phthalate	1300	U	4-Chlorophenyl phenyl ether	1300	U
Bis (2-ethylhexyl) phthalate	1300	U	Diethyl Phthalate	1300	U
1,2,4-Trichlorobenzene	1300	U	Hexachlorocyclopentadiene	1300	U

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis was conducted according to: Test Methods for Evaluating Solid Waste, SW-846 Method 8270C.

8270/625 layout

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Mr. Karl Kasper
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Portland ME 04102

May 31, 2006

CLIENT SAMPLE ID

Project Name: GP Old Town Mill

Project Number: 218736.002

Field Sample ID: SB-38 (0-4)

SAMPLE DATA

Lab Sample ID: 56637-16
Matrix: Solid
Percent Solid: 94
Dilution Factor: 5
Collection Date: 05/25/06
Lab Receipt Date: 05/27/06
Extraction Date: 05/30/06
Analysis Date: 05/31/06

ANALYTICAL RESULTS SEMI-VOLATILE ORGANICS

BASE NEUTRAL COMPOUND	Quantitation Limit µg/kg	Result µg/kg	BASE NEUTRAL COMPOUND	Quantitation Limit µg/kg	Result µg/kg
Acenaphthene	1300	**15300**	N-nitrosodimethylamine	1300	U
Acenaphthylene	1300	**16600**	N-nitroso-di-n-propylamine	1300	U
Anthracene	1300	**22000 E**	n-nitrosodiphenylamine	1300	**2060**
Benzo[a]anthracene	1300	U	Pyridine	1300	U
Benzo[a] pyrene	1300	**23400 E**	2-Methylnaphthalene	1300	**17200**
Benzo[b] fluoranthene	1300	**20300 E**	2-Chloronaphthalene	1300	U
Benzo[k] fluoranthene	1300	**21400 E**	Naphthalene	1300	**4500**
Benzo(g,h,i) perylene	1300	**7550**	Phenanthrene	1300	**76300 E**
Chrysene	1300	**28200 E**	Dibenzofuran	1300	**5350**
Dibenz [a,h] anthracene	1300	**1610**	Aniline	1300	U
Fluoranthene	1300	**2270**	4-Chloroaniline	1300	U
Fluorene	1300	**33800 E**	2-Nitroaniline	1300	U
Indeno [1,2,3-cd] pyrene	1300	**7080**	3-Nitroaniline	1300	U
Pyrene	1300	**122500 E**	4-Nitroaniline	1300	U
Hexachloroethane	1300	U	Carbazole	1300	U
Isophorone	1300	U			

Base Neutral Surrogate Standard Recovery

2-Fluorobiphenyl	69	%	d5-nitrobenzene	68	%	d14-p-terphenyl	185* %

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis was conducted according to: Test Methods for Evaluating Solid Waste, SW-846 Method 8270C.

COMMENTS: *Due to the reactive nature of this compound, the Benzidine quantitation limit is estimated. Results are expressed on a dry weight basis. Sample had low recovery of internal standard d12-Perylene. *Surrogate recovery outside laboratory acceptance criteria. Sample is being reanalyzed at a dilution.

8270/625 layout

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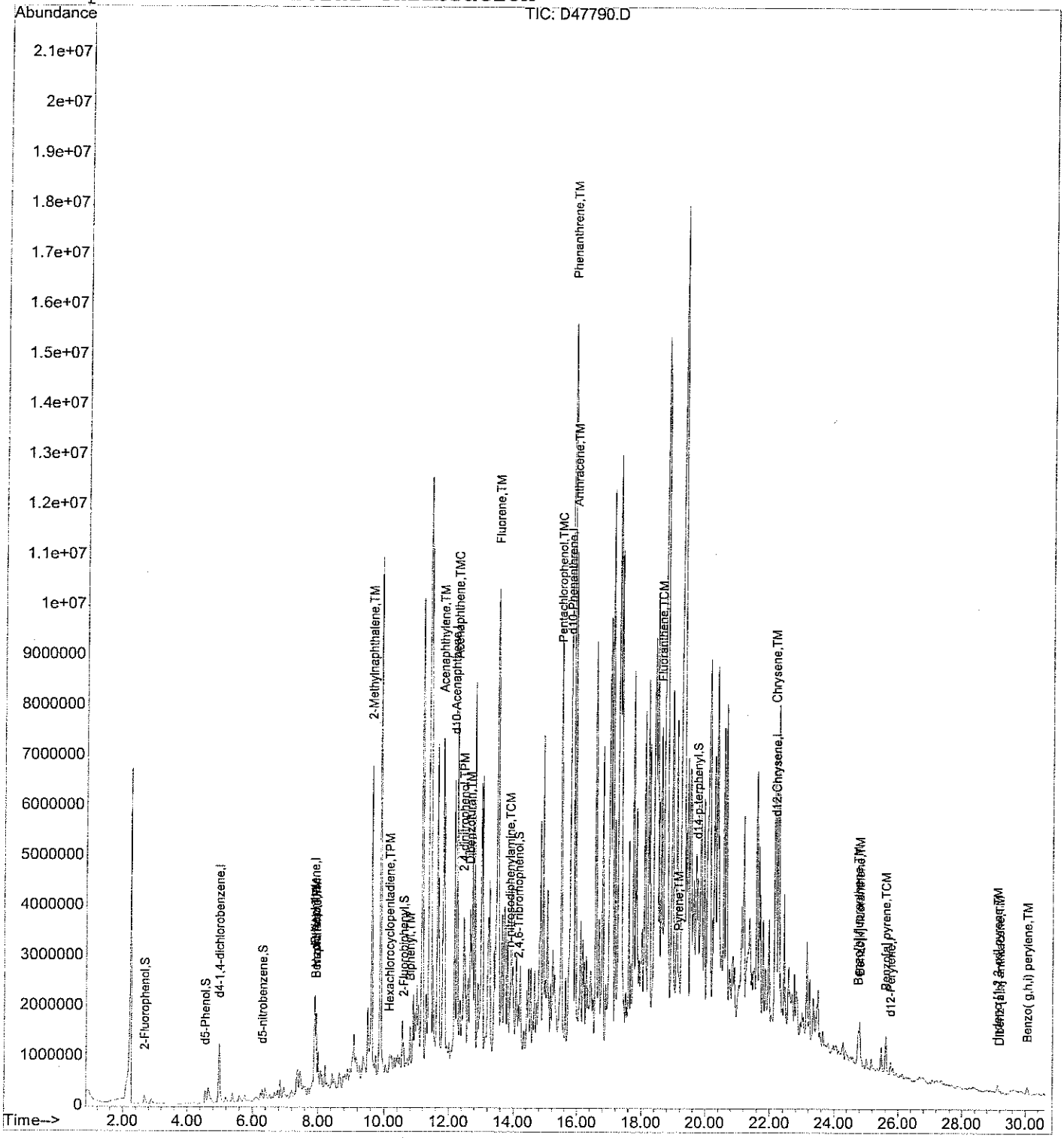
Quantitation Report

Data File : D:\HPCHEM\1\DATA\053006-D\D47790.D Vial: 20
Acq On : 31 May 2006 7:28 am Operator:
Sample : 56637-16,,1:5 Inst : D Instrum
Misc : SOIL Multiplr: 1.00
MS Integration Params: RTEINT.P
Quant Time: May 31 10:11 2006 Quant Results File: S05246.RES

Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Tue May 30 20:31:56 2006
Response via : Initial Calibration



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Portland ME 04102

June 1, 2006

CLIENT SAMPLE ID

Project Name: GP Old Town Mill

Project Number: 218736.002

Field Sample ID: SB-38 (0-4)

SAMPLE DATA

Lab Sample ID: 56637-16 DL
Matrix: Solid
Percent Solid: 94
Dilution Factor: 52
Collection Date: 05/25/06
Lab Receipt Date: 05/27/06
Extraction Date: 05/30/06
Analysis Date: 05/31/06

ANALYTICAL RESULTS SEMI-VOLATILE ORGANICS

ACID COMPOUND	Quantitation Limit µg/kg	Result µg/kg	ACID COMPOUND	Quantitation Limit µg/kg	Result µg/kg
2-Chlorophenol	13000	U	Pentachlorophenol	13000	U
4-Chloro-3-methylphenol	13000	U	Phenol	13000	U
2,4-Dichlorophenol	13000	U	2,4,5-Trichlorophenol	13000	U
2,4-Dimethylphenol	13000	U	2,4,6-Trichlorophenol	13000	U
2,4-dinitrophenol	13000	U	Benzoic Acid	13000	U
4,6-Dinitro-2-methylphenol	13000	U	2-Methylphenol	13000	U
2-Nitrophenol	13000	U	3+4-Methylphenol	13000	U
2,6-Dichlorophenol	13000	U	Benzyl Alcohol	13000	U
4-Nitrophenol	13000	U	2,3,4,6-Tetrachlorophenol	13000	U

Acid Surrogate Standard Recovery

2-Fluorophenol	* %	d5-Phenol	* %	2,4,6-Tribromophenol	* %

BASE NEUTRAL COMPOUND	Quantitation Limit µg/kg	Result µg/kg	BASE NEUTRAL COMPOUND	Quantitation Limit µg/kg	Result µg/kg
1,2-Dichlorobenzene	13000	U	Hexachlorobenzene	13000	U
1,3-Dichlorobenzene	13000	U	* Benzidine	13000	U
1,4-Dichlorobenzene	13000	U	3,3'-Dichlorobenzidine	13000	U
2,4-Dinitrotoluene	13000	U	Azobenzene	13000	U
2,6-Dinitrotoluene	13000	U	Bis(2-chloroethoxy)methane	13000	U
Nitrobenzene	13000	U	bis(2-chloroethyl) ether	13000	U
Hexachlorobutadiene	13000	U	bis(2-chloroisopropyl)ether	13000	U
Dimethyl Phthalate	13000	U	4-bromophenyl phenyl ether	13000	U
Di-n-butyl phthalate	13000	U	Butyl benzyl phthalate	13000	U
di-n-octyl-phthalate	13000	U	4-Chlorophenyl phenyl ether	13000	U
Bis (2-ethylhexyl) phthalate	13000	U	Diethyl Phthalate	13000	U
1,2,4-Trichlorobenzene	13000	U	Hexachlorocyclopentadiene	13000	U

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis was conducted according to: Test Methods for Evaluating Solid Waste, SW-846 Method 8270C.

Authorized signature Melina Sull

8270/625 layout

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June 1, 2006

CLIENT SAMPLE ID

Project Name: GP Old Town Mill

Project Number: 218736.002

Field Sample ID: SB-38 (0-4)

SAMPLE DATA

Lab Sample ID: 56637-16 DL
Matrix: Solid
Percent Solid: 94
Dilution Factor: 52
Collection Date: 05/25/06
Lab Receipt Date: 05/27/06
Extraction Date: 05/30/06
Analysis Date: 05/31/06

ANALYTICAL RESULTS SEMI-VOLATILE ORGANICS

BASE NEUTRAL COMPOUND	Quantitation Limit µg/kg	Result µg/kg	BASE NEUTRAL COMPOUND	Quantitation Limit µg/kg	Result µg/kg
Acenaphthene	13000	**14700**	N-nitrosodimethylamine	13000	U
Acenaphthylene	13000	**16200**	N-nitroso-di-n-propylamine	13000	U
Anthracene	13000	**21900**	n-nitrosodiphenylamine	13000	U
Benzo[a]anthracene	13000	**27400**	Pyridine	13000	U
Benzo[a] pyrene	13000	**22000**	2-Methylnaphthalene	13000	**14600**
Benzo[b] fluoranthene	13000	**14200**	2-Chloronaphthalene	13000	U
Benzo[k] fluoranthene	13000	**13300**	Naphthalene	13000	U
Benzo(g,h,i) perylene	13000	**12600 J**	Phenanthrene	13000	**100400**
Chrysene	13000	**29300**	Dibenzofuran	13000	U
Dibenz [a,h] anthracene	13000	U	Aniline	13000	U
Fluoranthene	13000	**68100**	4-Chloroaniline	13000	U
Fluorene	13000	**27500**	2-Nitroaniline	13000	U
Indeno [1,2,3-cd] pyrene	13000	**12300 J**	3-Nitroaniline	13000	U
Pyrene	13000	**80700**	4-Nitroaniline	13000	U
Hexachloroethane	13000	U	Carbazole	13000	U
Isophorone	13000	U			

Base Neutral Surrogate Standard Recovery

2-Fluorobiphenyl	* %	d5-nitrobenzene	* %	d14-p-terphenyl	* %

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis was conducted according to: Test Methods for Evaluating Solid Waste, SW-846 Method 8270C.

COMMENTS: *Due to the reactive nature of this compound, the Benzidine quantitation limit is estimated. Results are expressed on a dry weight basis. *The surrogates were diluted out.

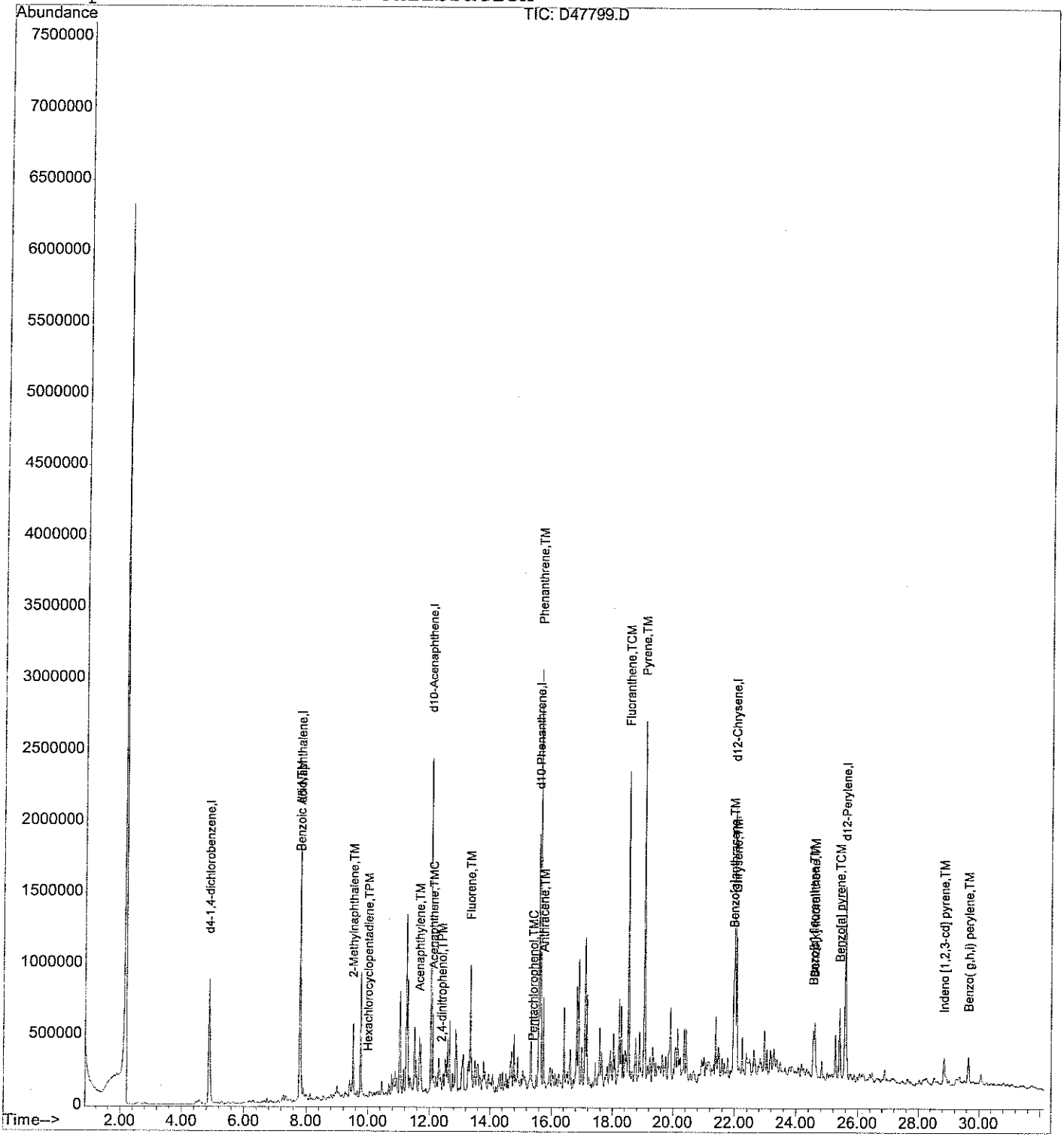
8270/625 layout

Authorized signature Melissa Tull

Quantitation Report

Data File : D:\HPCHEM\1\DATA\053106-D\D47799.D Vial: 6
Acq On : 31 May 2006 1:19 pm Operator:
Sample : 56637-16,,1:50 DL Inst : D Instrum
Misc : SOIL Multiplr: 1.00
MS Integration Params: RTEINT.P
Quant Time: May 31 13:52 2006 Quant Results File: S05246.RES

Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Wed May 31 12:09:53 2006
Response via : Initial Calibration



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May 31, 2006

SAMPLE DATA

CLIENT SAMPLE ID

Project Name:	GP Old Town Mill
Project Number:	218736.002
Field Sample ID:	SB-37 (0-4)

Lab Sample ID:	56637-17
Matrix:	Solid
Percent Solid:	93
Dilution Factor:	5
Collection Date:	05/25/06
Lab Receipt Date:	05/27/06
Extraction Date:	05/30/06
Analysis Date:	05/31/06

ANALYTICAL RESULTS SEMI-VOLATILE ORGANICS

ACID COMPOUND	Quantitation Limit µg/kg	Result µg/kg	ACID COMPOUND	Quantitation Limit µg/kg	Result µg/kg
2-Chlorophenol	1300	U	Pentachlorophenol	1300	U
4-Chloro-3-methylphenol	1300	U	Phenol	1300	U
2,4-Dichlorophenol	1300	U	2,4,5-Trichlorophenol	1300	U
2,4-Dimethylphenol	1300	U	2,4,6-Trichlorophenol	1300	U
2,4-dinitrophenol	1300	U	Benzoic Acid	1300	U
4,6-Dinitro-2-methylphenol	1300	U	2-Methylphenol	1300	U
2-Nitrophenol	1300	U	3+4-Methylphenol	1300	U
2,6-Dichlorophenol	1300	U	Benzyl Alcohol	1300	U
4-Nitrophenol	1300	U	2,3,4,6-Tetrachlorophenol	1300	U

Acid Surrogate Standard Recovery

2-Fluorophenol	57 %	d5-Phenol	61 %	2,4,6-Tribromophenol	69 %

BASE NEUTRAL COMPOUND	Quantitation Limit µg/kg	Result µg/kg	BASE NEUTRAL COMPOUND	Quantitation Limit µg/kg	Result µg/kg
1,2-Dichlorobenzene	1300	U	Hexachlorobenzene	1300	U
1,3-Dichlorobenzene	1300	U	* Benzidine	1300	U
1,4-Dichlorobenzene	1300	U	3,3'-Dichlorobenzidine	1300	U
2,4-Dinitrotoluene	1300	U	Azobenzene	1300	U
2,6-Dinitrotoluene	1300	U	Bis(2-chloroethoxy)methane	1300	U
Nitrobenzene	1300	U	bis(2-chloroethyl) ether	1300	U
Hexachlorobutadiene	1300	U	bis(2-chloroisopropyl)ether	1300	U
Dimethyl Phthalate	1300	U	4-bromophenyl phenyl ether	1300	U
Di-n-butyl phthalate	1300	U	Butyl benzyl phthalate	1300	U
di-n-octyl-phthalate	1300	U	4-Chlorophenyl phenyl ether	1300	U
Bis (2-ethylhexyl) phthalate	1300	U	Diethyl Phthalate	1300	U
1,2,4-Trichlorobenzene	1300	U	Hexachlorocyclopentadiene	1300	U

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis was conducted according to: Test Methods for Evaluating Solid Waste, SW-846 Method 8270C.

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8270/625 layout

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May 31, 2006

CLIENT SAMPLE ID

Project Name: GP Old Town Mill

Project Number: 218736.002

Field Sample ID: SB-37 (0-4)

SAMPLE DATA

Lab Sample ID: 56637-17
Matrix: Solid
Percent Solid: 93
Dilution Factor: 5
Collection Date: 05/25/06
Lab Receipt Date: 05/27/06
Extraction Date: 05/30/06
Analysis Date: 05/31/06

ANALYTICAL RESULTS SEMI-VOLATILE ORGANICS

BASE NEUTRAL COMPOUND	Quantitation Limit µg/kg	Result µg/kg	BASE NEUTRAL COMPOUND	Quantitation Limit µg/kg	Result µg/kg
Acenaphthene	1300	U	N-nitrosodimethylamine	1300	U
Acenaphthylene	1300	**9030**	N-nitroso-di-n-propylamine	1300	U
Anthracene	1300	**2230**	n-nitrosodiphenylamine	1300	U
Benzo[a]anthracene	1300	**12800**	Pyridine	1300	U
Benzo[a] pyrene	1300	**17000**	2-Methylnaphthalene	1300	**671 J**
Benzo[b] fluoranthene	1300	**15600**	2-Chloronaphthalene	1300	U
Benzo[k] fluoranthene	1300	**12900**	Naphthalene	1300	**680 J**
Benzo(g,h,i) perylene	1300	**7240**	Phenanthrene	1300	**8780**
Chrysene	1300	**16600**	Dibenzofuran	1300	U
Dibenz [a,h] anthracene	1300	**1500**	Aniline	1300	U
Fluoranthene	1300	U	4-Chloroaniline	1300	U
Fluorene	1300	**718 J**	2-Nitroaniline	1300	U
Indeno [1,2,3-cd] pyrene	1300	**6690**	3-Nitroaniline	1300	U
Pyrene	1300	**92900 E**	4-Nitroaniline	1300	U
Hexachloroethane	1300	U	Carbazole	1300	U
Isophorone	1300	U			

Base Neutral Surrogate Standard Recovery

2-Fluorobiphenyl	67	%	d5-nitrobenzene	59	%	d14-p-terphenyl	166* %

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis was conducted according to: Test Methods for Evaluating Solid Waste, SW-846 Method 8270C.

COMMENTS: *Due to the reactive nature of this compound, the Benzidine quantitation limit is estimated.
Results are expressed on a dry weight basis. Sample had low recovery of internal standards d12-Chrysene and d12-Perylene.
*Surrogate recovery outside laboratory acceptance criteria. Sample is being reanalyzed at a dilution.

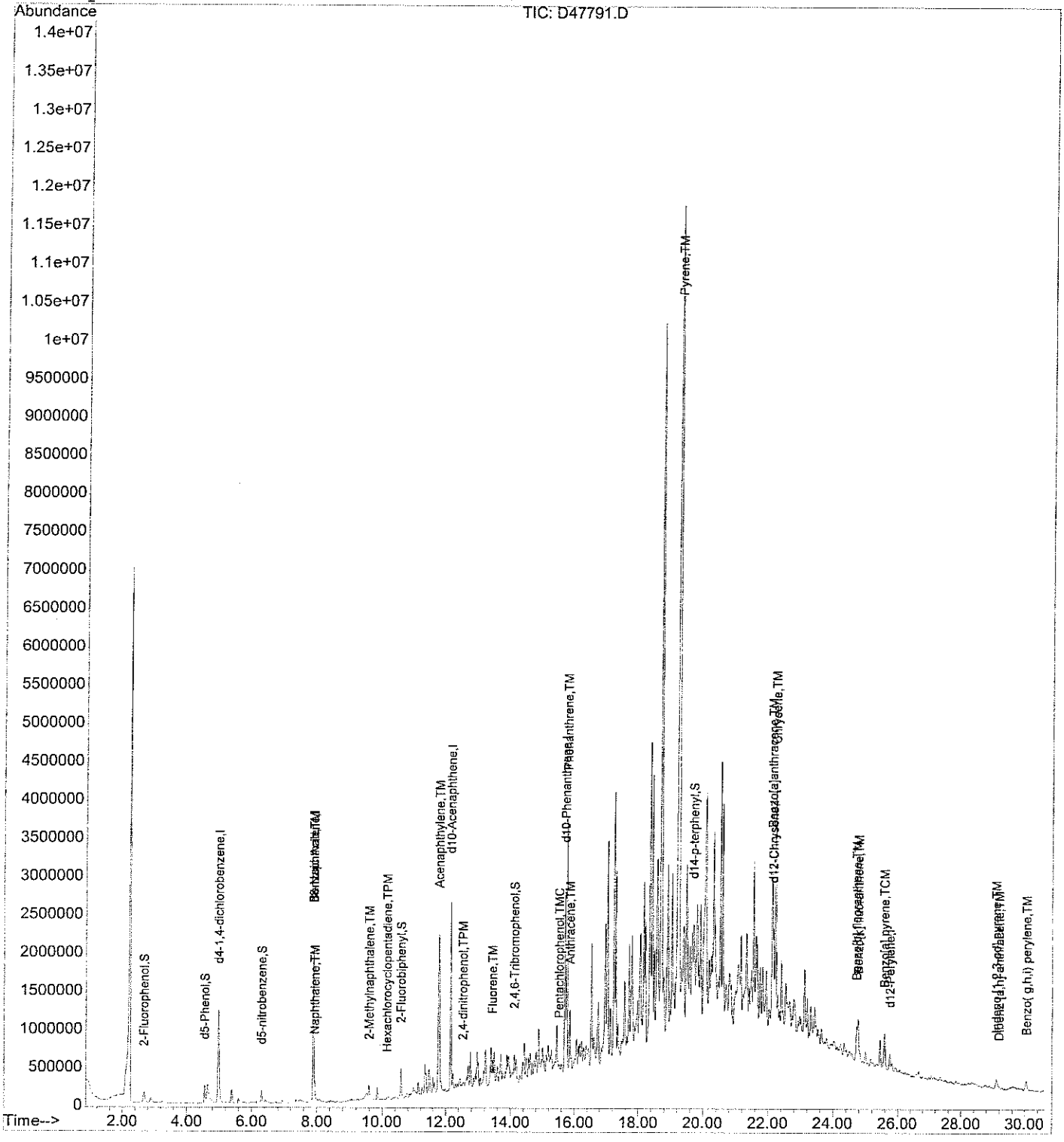
8270/625 layout

Authorized signature

Quantitation Report

Data File : D:\HPCHEM\1\DATA\053006-D\D47791.D Vial: 21
Acq On : 31 May 2006 8:01 am Operator:
Sample : 56637-17,,1:5 Inst : D Instrum
Misc : SOIL Multiplr: 1.00
MS Integration Params: RTEINT.P
Quant Time: May 31 10:41 2006 Quant Results File: S05246.RES

Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Tue May 30 20:31:56 2006
Response via : Initial Calibration



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Mr. Karl Kasper
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41 Hutchins Drive
Portland ME 04102

June 1, 2006

CLIENT SAMPLE ID

Project Name: GP Old Town Mill

Project Number: 218736.002

Field Sample ID: SB-37 (0-4)

SAMPLE DATA

Lab Sample ID: 56637-17 DL
Matrix: Solid
Percent Solid: 93
Dilution Factor: 21
Collection Date: 05/25/06
Lab Receipt Date: 05/27/06
Extraction Date: 05/30/06
Analysis Date: 05/31/06

ANALYTICAL RESULTS SEMI-VOLATILE ORGANICS

ACID COMPOUND	Quantitation Limit µg/kg	Result µg/kg	ACID COMPOUND	Quantitation Limit µg/kg	Result µg/kg
2-Chlorophenol	5300	U	Pentachlorophenol	5300	U
4-Chloro-3-methylphenol	5300	U	Phenol	5300	U
2,4-Dichlorophenol	5300	U	2,4,5-Trichlorophenol	5300	U
2,4-Dimethylphenol	5300	U	2,4,6-Trichlorophenol	5300	U
2,4-dinitrophenol	5300	U	Benzoic Acid	5300	U
4,6-Dinitro-2-methylphenol	5300	U	2-Methylphenol	5300	U
2-Nitrophenol	5300	U	3+4-Methylphenol	5300	U
2,6-Dichlorophenol	5300	U	Benzyl Alcohol	5300	U
4-Nitrophenol	5300	U	2,3,4,6-Tetrachlorophenol	5300	U

Acid Surrogate Standard Recovery

2-Fluorophenol	* %	d5-Phenol	* %	2,4,6-Tribromophenol	* %

BASE NEUTRAL COMPOUND	Quantitation Limit µg/kg	Result µg/kg	BASE NEUTRAL COMPOUND	Quantitation Limit µg/kg	Result µg/kg
1,2-Dichlorobenzene	5300	U	Hexachlorobenzene	5300	U
1,3-Dichlorobenzene	5300	U	* Benzidine	5300	U
1,4-Dichlorobenzene	5300	U	3,3'-Dichlorobenzidine	5300	U
2,4-Dinitrotoluene	5300	U	Azobenzene	5300	U
2,6-Dinitrotoluene	5300	U	Bis(2-chloroethoxy)methane	5300	U
Nitrobenzene	5300	U	bis(2-chloroethyl) ether	5300	U
Hexachlorobutadiene	5300	U	bis(2-chloroisopropyl)ether	5300	U
Dimethyl Phthalate	5300	U	4-bromophenyl phenyl ether	5300	U
Di-n-butyl phthalate	5300	U	Butyl benzyl phthalate	5300	U
di-n-octyl-phthalate	5300	U	4-Chlorophenyl phenyl ether	5300	U
Bis (2-ethylhexyl) phthalate	5300	U	Diethyl Phthalate	5300	U
1,2,4-Trichlorobenzene	5300	U	Hexachlorocyclopentadiene	5300	U

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis was conducted according to: Test Methods for Evaluating Solid Waste, SW-846 Method 8270C.

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8270/625 layout

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Portland ME 04102

June 1, 2006

CLIENT SAMPLE ID

Project Name: GP Old Town Mill

Project Number: 218736.002

Field Sample ID: SB-37 (0-4)

SAMPLE DATA

Lab Sample ID: 56637-17 DL
Matrix: Solid
Percent Solid: 93
Dilution Factor: 21
Collection Date: 05/25/06
Lab Receipt Date: 05/27/06
Extraction Date: 05/30/06
Analysis Date: 05/31/06

ANALYTICAL RESULTS SEMI-VOLATILE ORGANICS

BASE NEUTRAL COMPOUND	Quantitation Limit µg/kg	Result µg/kg	BASE NEUTRAL COMPOUND	Quantitation Limit µg/kg	Result µg/kg
Acenaphthene	5300	U	N-nitrosodimethylamine	5300	U
Acenaphthylene	5300	**10800**	N-nitroso-di-n-propylamine	5300	U
Anthracene	5300	**2890 J**	n-nitrosodiphenylamine	5300	U
Benzo[a]anthracene	5300	**13800**	Pyridine	5300	U
Benzo[a] pyrene	5300	**18600**	2-Methylnaphthalene	5300	U
Benzo[b] fluoranthene	5300	**16700**	2-Chloronaphthalene	5300	U
Benzo[k] fluoranthene	5300	**12900**	Naphthalene	5300	U
Benzo(g,h,i) perylene	5300	**4080 J**	Phenanthrene	5300	**9220**
Chrysene	5300	**18200**	Dibenzofuran	5300	U
Dibenz [a,h] anthracene	5300	U	Aniline	5300	U
Fluoranthene	5300	**26800**	4-Chloroaniline	5300	U
Fluorene	5300	U	2-Nitroaniline	5300	U
Indeno [1,2,3-cd] pyrene	5300	**4730 J**	3-Nitroaniline	5300	U
Pyrene	5300	**39100**	4-Nitroaniline	5300	U
Hexachloroethane	5300	U	Carbazole	5300	U
Isophorone	5300	U			

Base Neutral Surrogate Standard Recovery

2-Fluorobiphenyl	* %	d5-nitrobenzene	* %	d14-p-terphenyl	* %

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis was conducted according to: Test Methods for Evaluating Solid Waste, SW-846 Method 8270C.

COMMENTS: *Due to the reactive nature of this compound, the Benzidine quantitation limit is estimated. Results are expressed on a dry weight basis. *The surrogates were diluted out.

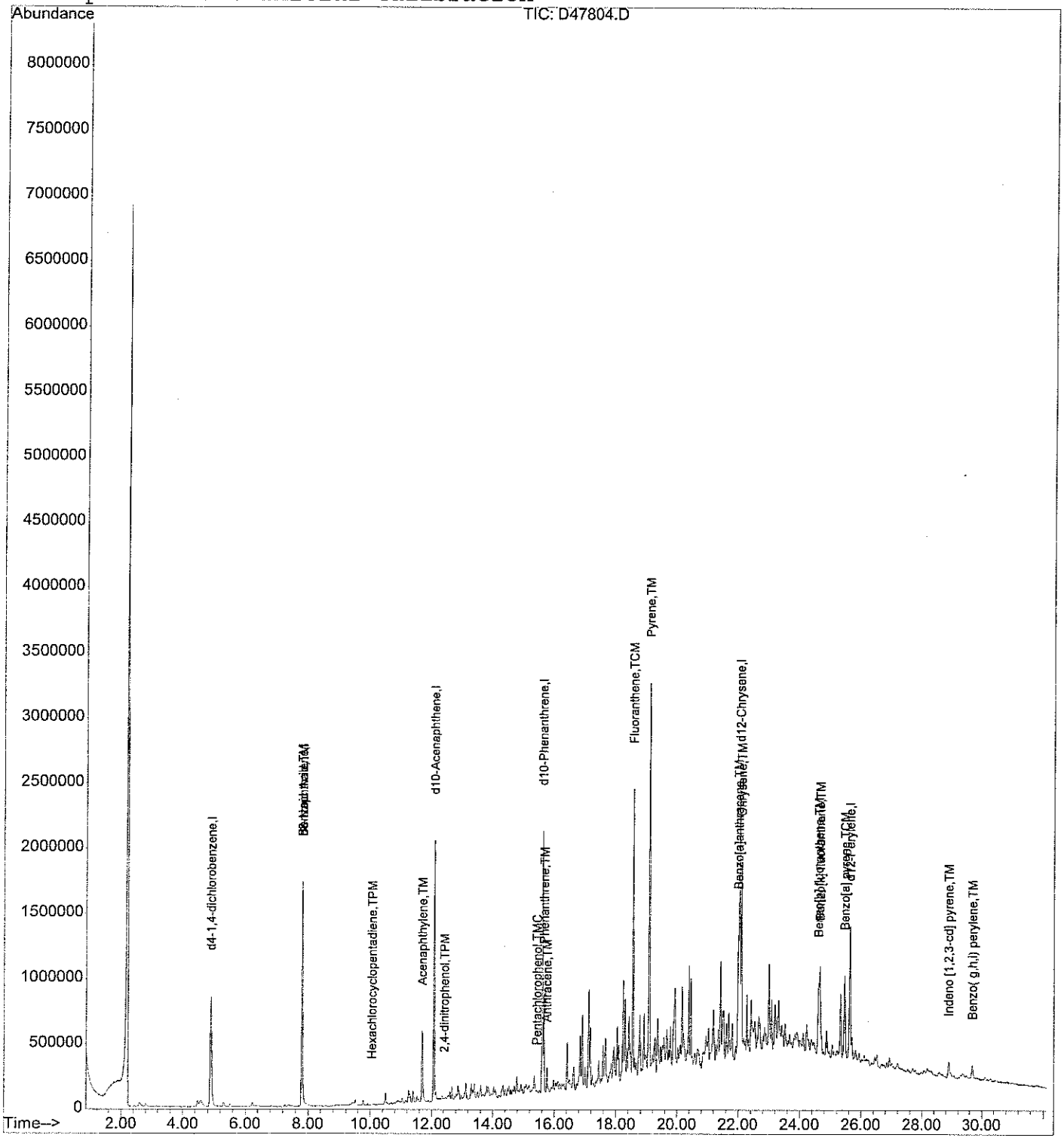
8270/625 layout

Authorized signature Mylena Tulli

Quantitation Report

Data File : D:\HPCHEM\1\DATA\053106-D\D47804.D Vial: 11
Acq On : 31 May 2006 4:43 pm Operator:
Sample : 56637-17,,1:20 Inst : D Instrum
Misc : SOIL Multiplr: 1.00
MS Integration Params: RTEINT.P
Quant Time: May 31 20:11 2006 Quant Results File: S05246.RES

Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Wed May 31 19:58:02 2006
Response via : Initial Calibration



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May 31, 2006

CLIENT SAMPLE ID

Project Name: GP Old Town Mill

Project Number: 218736.002

Field Sample ID: SB-15 (0-4)

SAMPLE DATA

Lab Sample ID: 56637-18
Matrix: Solid
Percent Solid: 88
Dilution Factor: 1.1
Collection Date: 05/25/06
Lab Receipt Date: 05/27/06
Extraction Date: 05/30/06
Analysis Date: 05/31/06

ANALYTICAL RESULTS SEMI-VOLATILE ORGANICS

ACID COMPOUND	Quantitation Limit µg/kg	Result µg/kg	ACID COMPOUND	Quantitation Limit µg/kg	Result µg/kg
2-Chlorophenol	280	U	Pentachlorophenol	280	U
4-Chloro-3-methylphenol	280	U	Phenol	280	U
2,4-Dichlorophenol	280	U	2,4,5-Trichlorophenol	280	U
2,4-Dimethylphenol	280	U	2,4,6-Trichlorophenol	280	U
2,4-dinitrophenol	280	U	Benzoic Acid	280	U
4,6-Dinitro-2-methylphenol	280	U	2-Methylphenol	280	U
2-Nitrophenol	280	U	3+4-Methylphenol	280	U
2,6-Dichlorophenol	280	U	Benzyl Alcohol	280	U
4-Nitrophenol	280	U	2,3,4,6-Tetrachlorophenol	280	U

Acid Surrogate Standard Recovery

2-Fluorophenol	79 %	d5-Phenol	89 %	2,4,6-Tribromophenol	94 %

BASE NEUTRAL COMPOUND	Quantitation Limit µg/kg	Result µg/kg	BASE NEUTRAL COMPOUND	Quantitation Limit µg/kg	Result µg/kg
1,2-Dichlorobenzene	280	U	Hexachlorobenzene	280	U
1,3-Dichlorobenzene	280	U	* Benzidine	280	U
1,4-Dichlorobenzene	280	U	3,3'-Dichlorobenzidine	280	U
2,4-Dinitrotoluene	280	U	Azobenzene	280	U
2,6-Dinitrotoluene	280	U	Bis(2-chloroethoxy)methane	280	U
Nitrobenzene	280	U	bis(2-chloroethyl) ether	280	U
Hexachlorobutadiene	280	U	bis(2-chloroisopropyl)ether	280	U
Dimethyl Phthalate	280	U	4-bromophenyl phenyl ether	280	U
Di-n-butyl phthalate	280	U	Butyl benzyl phthalate	280	U
di-n-octyl-phthalate	280	U	4-Chlorophenyl phenyl ether	280	U
Bis (2-ethylhexyl) phthalate	280	U	Diethyl Phthalate	280	U
1,2,4-Trichlorobenzene	280	U	Hexachlorocyclopentadiene	280	U

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis was conducted according to: Test Methods for Evaluating Solid Waste, SW-846 Method 8270C.

Authorized signature

8270/625 layout

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May 31, 2006

CLIENT SAMPLE ID

Project Name: GP Old Town Mill

Project Number: 218736.002

Field Sample ID: SB-15 (0-4)

SAMPLE DATA

Lab Sample ID: 56637-18
Matrix: Solid
Percent Solid: 88
Dilution Factor: 1.1
Collection Date: 05/25/06
Lab Receipt Date: 05/27/06
Extraction Date: 05/30/06
Analysis Date: 05/31/06

ANALYTICAL RESULTS SEMI-VOLATILE ORGANICS

BASE NEUTRAL COMPOUND	Quantitation Limit µg/kg	Result µg/kg	BASE NEUTRAL COMPOUND	Quantitation Limit µg/kg	Result µg/kg
Acenaphthene	280	U	N-nitrosodimethylamine	280	U
Acenaphthylene	280	U	N-nitroso-di-n-propylamine	280	U
Anthracene	280	U	n-nitrosodiphenylamine	280	U
Benzo[a]anthracene	280	U	Pyridine	280	U
Benzo[a] pyrene	280	U	2-Methylnaphthalene	280	U
Benzo[b] fluoranthene	280	U	2-Chloronaphthalene	280	U
Benzo[k] fluoranthene	280	U	Naphthalene	280	U
Benzo(g,h,i) perylene	280	U	Phenanthrene	280	U
Chrysene	280	U	Dibenzofuran	280	U
Dibenz [a,h] anthracene	280	U	Aniline	280	U
Fluoranthene	280	U	4-Chloroaniline	280	U
Fluorene	280	U	2-Nitroaniline	280	U
Indeno [1,2,3-cd] pyrene	280	U	3-Nitroaniline	280	U
Pyrene	280	U	4-Nitroaniline	280	U
Hexachloroethane	280	U	Carbazole	280	U
Isophorone	280	U			

Base Neutral Surrogate Standard Recovery

2-Fluorobiphenyl	79	%	d5-nitrobenzene	87	%	d14-p-terphenyl	91 %

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis was conducted according to: Test Methods for Evaluating Solid Waste, SW-846 Method 8270C.

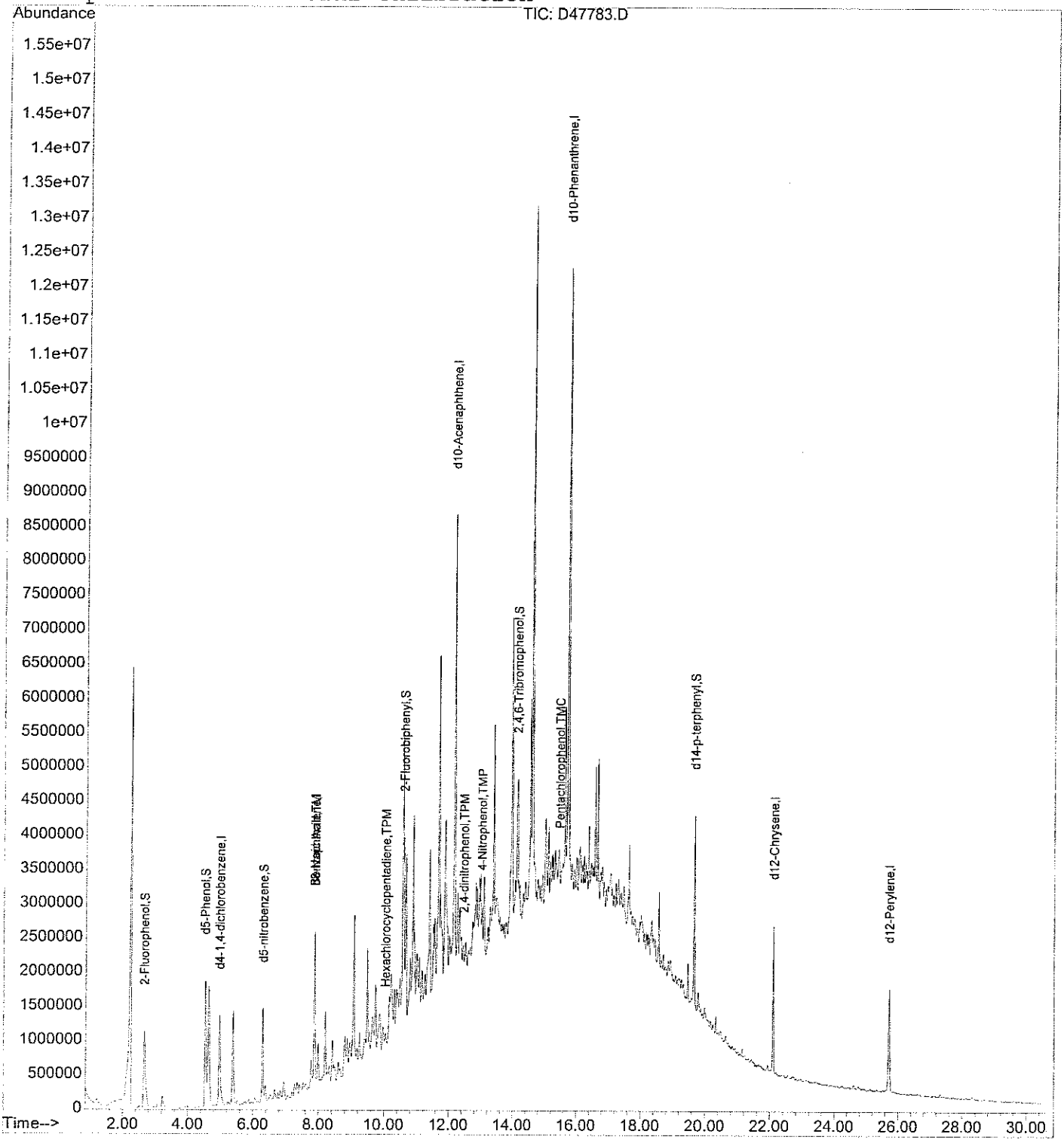
COMMENTS: *Due to the reactive nature of this compound, the Benzidine quantitation limit is estimated.
Results are expressed on a dry weight basis.

8270/625 layout

Authorized signature

Data File : D:\HPCHEM\1\DATA\053006-D\D47783.D Vial: 13
Acq On : 31 May 2006 3:37 am Operator:
Sample : 56637-18 Inst : D Instrum
Misc : SOIL Multiplr: 1.00
MS Integration Params: RTEINT.P
Quant Time: May 31 8:21 2006 Quant Results File: S05246.RES

Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Tue May 30 20:31:56 2006
Response via : Initial Calibration



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May 31, 2006

CLIENT SAMPLE ID

Project Name: GP Old Town Mill

Project Number: 218736.002

Field Sample ID: SB-41 (0-4)

SAMPLE DATA

Lab Sample ID: 56637-19
Matrix: Solid
Percent Solid: 91
Dilution Factor: 1.1
Collection Date: 05/26/06
Lab Receipt Date: 05/27/06
Extraction Date: 05/30/06
Analysis Date: 05/31/06

ANALYTICAL RESULTS SEMI-VOLATILE ORGANICS

ACID COMPOUND	Quantitation Limit µg/kg	Result µg/kg	ACID COMPOUND	Quantitation Limit µg/kg	Result µg/kg
2-Chlorophenol	280	U	Pentachlorophenol	280	U
4-Chloro-3-methylphenol	280	U	Phenol	280	U
2,4-Dichlorophenol	280	U	2,4,5-Trichlorophenol	280	U
2,4-Dimethylphenol	280	U	2,4,6-Trichlorophenol	280	U
2,4-dinitrophenol	280	U	Benzoic Acid	280	U
4,6-Dinitro-2-methylphenol	280	U	2-Methylphenol	280	U
2-Nitrophenol	280	U	3+4-Methylphenol	280	U
2,6-Dichlorophenol	280	U	Benzyl Alcohol	280	U
4-Nitrophenol	280	U	2,3,4,6-Tetrachlorophenol	280	U

Acid Surrogate Standard Recovery

2-Fluorophenol	73 %	d5-Phenol	84 %	2,4,6-Tribromophenol	95 %

BASE NEUTRAL COMPOUND	Quantitation Limit µg/kg	Result µg/kg	BASE NEUTRAL COMPOUND	Quantitation Limit µg/kg	Result µg/kg
1,2-Dichlorobenzene	280	U	Hexachlorobenzene	280	U
1,3-Dichlorobenzene	280	U	* Benzidine	280	U
1,4-Dichlorobenzene	280	U	3,3'-Dichlorobenzidine	280	U
2,4-Dinitrotoluene	280	U	Azobenzene	280	U
2,6-Dinitrotoluene	280	U	Bis(2-chloroethoxy)methane	280	U
Nitrobenzene	280	U	bis(2-chloroethyl) ether	280	U
Hexachlorobutadiene	280	U	bis(2-chloroisopropyl)ether	280	U
Dimethyl Phthalate	280	U	4-bromophenyl phenyl ether	280	U
Di-n-butyl phthalate	280	U	Butyl benzyl phthalate	280	U
di-n-octyl-phthalate	280	U	4-Chlorophenyl phenyl ether	280	U
Bis (2-ethylhexyl) phthalate	280	U	Diethyl Phthalate	280	U
1,2,4-Trichlorobenzene	280	U	Hexachlorocyclopentadiene	280	U

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis was conducted according to: Test Methods for Evaluating Solid Waste, SW-846 Method 8270C.

Authorized signature Melissa Hall

8270/625 layout

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May 31, 2006

CLIENT SAMPLE ID

Project Name: GP Old Town Mill

Project Number: 218736.002

Field Sample ID: SB-41 (0-4)

SAMPLE DATA

Lab Sample ID: 56637-19
Matrix: Solid
Percent Solid: 91
Dilution Factor: 1.1
Collection Date: 05/26/06
Lab Receipt Date: 05/27/06
Extraction Date: 05/30/06
Analysis Date: 05/31/06

ANALYTICAL RESULTS SEMI-VOLATILE ORGANICS

BASE NEUTRAL COMPOUND	Quantitation Limit µg/kg	Result µg/kg	BASE NEUTRAL COMPOUND	Quantitation Limit µg/kg	Result µg/kg
Acenaphthene	280	U	N-nitrosodimethylamine	280	U
Acenaphthylene	280	U	N-nitroso-di-n-propylamine	280	U
Anthracene	280	U	n-nitrosodiphenylamine	280	U
Benzo[a]anthracene	280	U	Pyridine	280	U
Benzo[a] pyrene	280	U	2-Methylnaphthalene	280	U
Benzo[b] fluoranthene	280	U	2-Chloronaphthalene	280	U
Benzo[k] fluoranthene	280	U	Naphthalene	280	U
Benzo(g,h,i) perylene	280	U	Phenanthrene	280	U
Chrysene	280	U	Dibenzofuran	280	U
Dibenz [a,h] anthracene	280	U	Aniline	280	U
Fluoranthene	280	**158 J**	4-Chloroaniline	280	U
Fluorene	280	U	2-Nitroaniline	280	U
Indeno [1,2,3-cd] pyrene	280	U	3-Nitroaniline	280	U
Pyrene	280	U	4-Nitroaniline	280	U
Hexachloroethane	280	U	Carbazole	280	U
Isophorone	280	U			

Base Neutral Surrogate Standard Recovery

2-Fluorobiphenyl	83	%	d5-nitrobenzene	77	%	d14-p-terphenyl	70 %

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis was conducted according to: Test Methods for Evaluating Solid Waste, SW-846 Method 8270C.

COMMENTS: *Due to the reactive nature of this compound, the Benzidine quantitation limit is estimated.
Results are expressed on a dry weight basis.

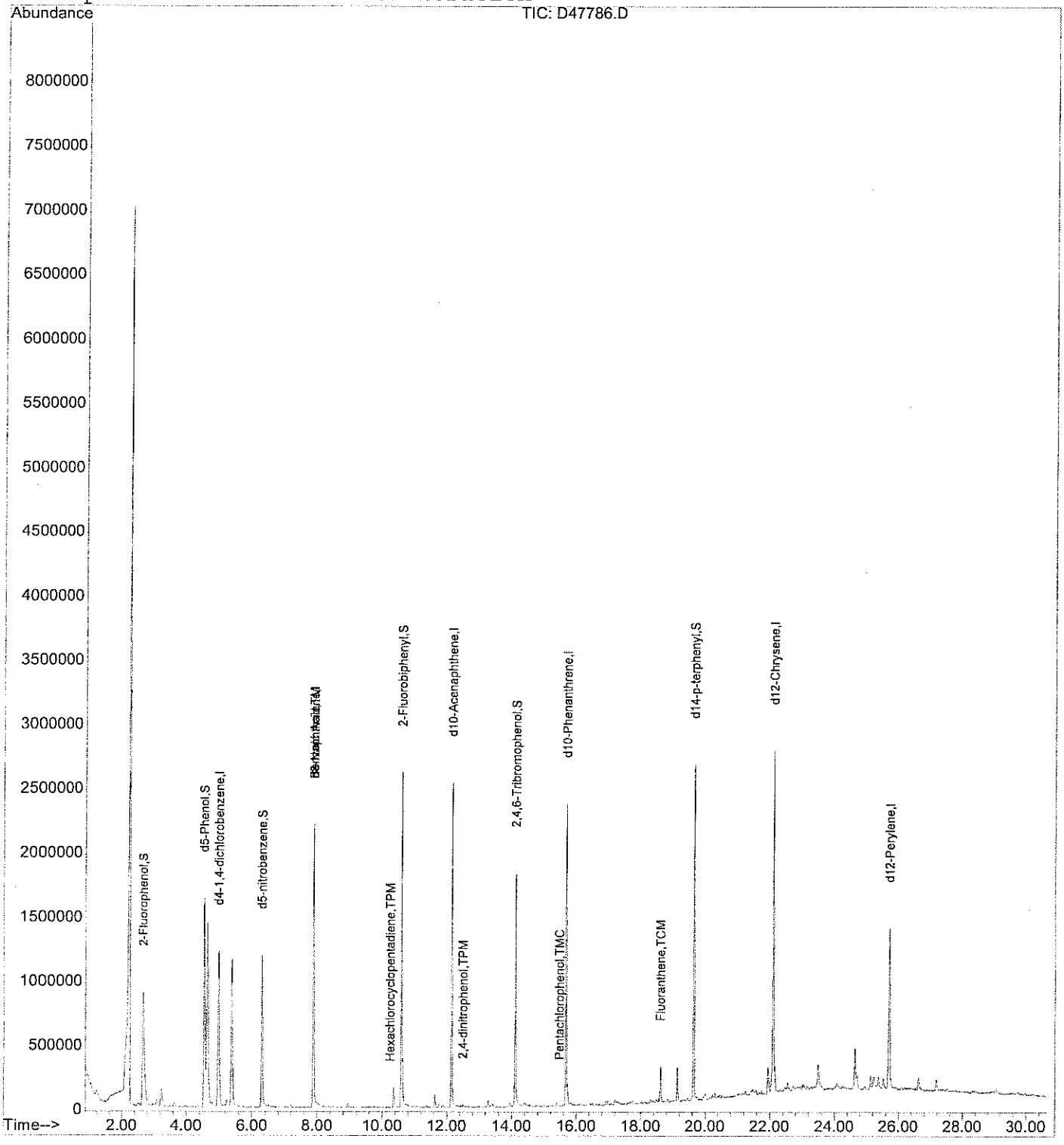
8270/625 layout

Authorized signature

Quantitation Report

Data File : D:\HPCHEM\1\DATA\053006-D\D47786.D Vial: 16
Acq On : 31 May 2006 5:16 am Operator:
Sample : 56637-19 Inst : D Instrum
Misc : SOIL Multiplr: 1.00
MS Integration Params: RTEINT.P
Quant Time: May 31 8:21 2006 Quant Results File: S05246.RES

Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Tue May 30 20:31:56 2006
Response via : Initial Calibration



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May 31, 2006

CLIENT SAMPLE ID

Project Name: GP Old Town Mill

Project Number: 218736.002

Field Sample ID: SB-15

SAMPLE DATA

Lab Sample ID: 56637-20
Matrix: Aqueous
Percent Solid: N/A
Dilution Factor: 1.0
Collection Date: 05/26/06
Lab Receipt Date: 05/27/06
Extraction Date: 05/30/06
Analysis Date: 05/30/06

ANALYTICAL RESULTS SEMI-VOLATILE ORGANICS

ACID COMPOUND	Quantitation Limit µg/L	Result µg/L	ACID COMPOUND	Quantitation Limit µg/L	Result µg/L
2-Chlorophenol	5	U	Pentachlorophenol	10	U
4-Chloro-3-methylphenol	10	U	Phenol	5	U
2,4-Dichlorophenol	5	U	2,4,5-Trichlorophenol	5	U
2,4-Dimethylphenol	5	U	2,4,6-Trichlorophenol	5	U
2,4-dinitrophenol	5	U	Benzoic Acid	10	U
4,6-Dinitro-2-methylphenol	5	U	2-Methylphenol	5	U
2-Nitrophenol	5	U	3+4-Methylphenol	5	U
2,6-Dichlorophenol	5	U	Benzyl Alcohol	5	U
4-Nitrophenol	5	U	2,3,4,6-Tetrachlorophenol	5	U

Acid Surrogate Standard Recovery

2-Fluorophenol	2 * %	d5-Phenol	1 * %	2,4,6-Tribromophenol	5* %

BASE NEUTRAL COMPOUND	Quantitation Limit µg/L	Result µg/L	BASE NEUTRAL COMPOUND	Quantitation Limit µg/L	Result µg/L
1,2-Dichlorobenzene	2	U	Hexachlorobenzene	2	U
1,3-Dichlorobenzene	2	U	* Benzidine	20	U
1,4-Dichlorobenzene	2	U	3,3'-Dichlorobenzidine	20	U
2,4-Dinitrotoluene	2	U	Azobenzene	2	U
2,6-Dinitrotoluene	2	U	Bis(2-chloroethoxy)methane	2	U
Nitrobenzene	2	U	bis(2-chloroethyl) ether	2	U
Hexachlorobutadiene	2	U	bis(2-chloroisopropyl)ether	2	U
Dimethyl Phthalate	2	U	4-bromophenyl phenyl ether	2	U
Di-n-butyl phthalate	2	U	Butyl benzyl phthalate	2	U
di-n-octyl-phthalate	2	U	4-Chlorophenyl phenyl ether	2	U
Bis (2-ethylhexyl) phthalate	2	U	Diethyl Phthalate	2	U
1,2,4-Trichlorobenzene	2	U	Hexachlorocyclopentadiene	2	U

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis was conducted according to: Test Methods for Evaluating Solid Waste, SW-846 Method 8270C.

Authorized signature

8270/625 layout

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May 31, 2006

CLIENT SAMPLE ID

Project Name: GP Old Town Mill

Project Number: 218736.002

Field Sample ID: SB-15

SAMPLE DATA

Lab Sample ID: 56637-20
Matrix: Aqueous
Percent Solid: N/A
Dilution Factor: 1.0
Collection Date: 05/26/06
Lab Receipt Date: 05/27/06
Extraction Date: 05/30/06
Analysis Date: 05/30/06

ANALYTICAL RESULTS SEMI-VOLATILE ORGANICS

BASE NEUTRAL COMPOUND	Quantitation Limit µg/L	Result µg/L	BASE NEUTRAL COMPOUND	Quantitation Limit µg/L	Result µg/L
Acenaphthene	2	U	N-nitrosodimethylamine	2	U
Acenaphthylene	2	U	N-nitroso-di-n-propylamine	2	U
Anthracene	2	U	n-nitrosodiphenylamine	2	U
Benzo[a]anthracene	2	U	Pyridine	2	U
Benzo[a] pyrene	2	U	2-Methylnaphthalene	2	U
Benzo[b] fluoranthene	2	U	2-Chloronaphthalene	2	U
Benzo[k] fluoranthene	2	U	Naphthalene	2	U
Benzo(g,h,i) perylene	2	U	Phenanthrene	2	U
Chrysene	2	U	Dibenzofuran	2	U
Dibenz [a,h] anthracene	2	U	Aniline	2	U
Fluoranthene	2	U	4-Chloroaniline	2	U
Fluorene	2	U	2-Nitroaniline	2	U
Indeno [1,2,3-cd] pyrene	2	U	3-Nitroaniline	2	U
Pyrene	2	U	4-Nitroaniline	2	U
Hexachloroethane	2	U	Carbazole	2	U
Isophorone	2	U			

Base Neutral Surrogate Standard Recovery

2-Fluorobiphenyl	72	%	d5-nitrobenzene	70	%	d14-p-terphenyl	73 %

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis was conducted according to: Test Methods for Evaluating Solid Waste, SW-846 Method 8270C.

COMMENTS: *Due to the reactive nature of this compound, the Benzidine quantitation limit is estimated.
*Surrogate recovery outside laboratory acceptance criteria. Sample was reanalyzed with similar results.

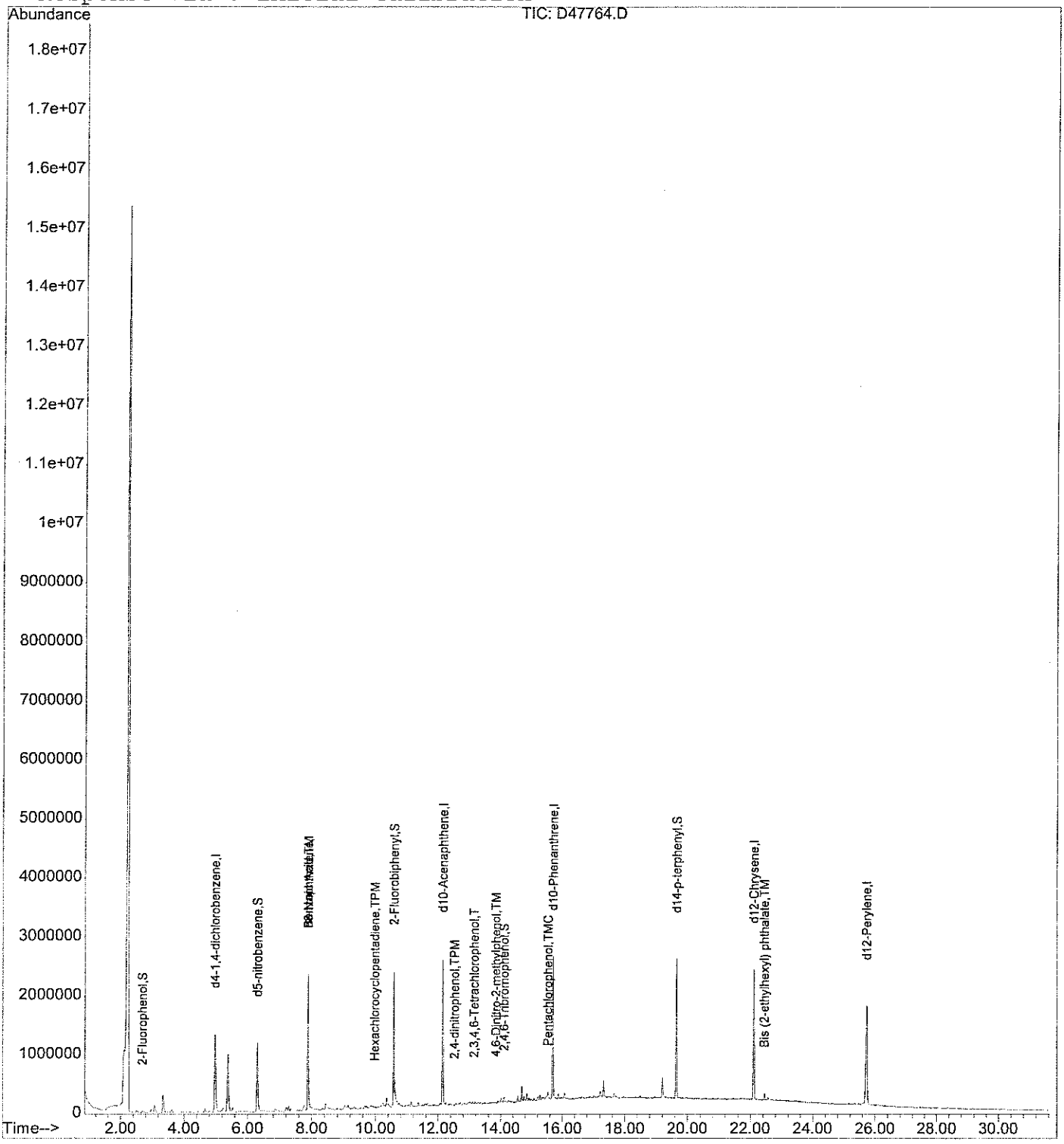
8270/625 layout

Authorized signature

Quantitation Report

Data File : D:\HPCHEM\1\DATA\053006-D\D47764.D Vial: 10
Acq On : 30 May 2006 1:08 pm Operator:
Sample : 56637-20 Inst : D Instrum
Misc : Multiplr: 1.00
MS Integration Params: RTEINT.P
Quant Time: May 30 13:19 2006 Quant Results File: S05246.RES

Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Tue May 30 09:52:41 2006
Response via : Initial Calibration



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Mr. Karl Kasper
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May 31, 2006

CLIENT SAMPLE ID

Project Name: GP Old Town Mill

Project Number: 218736.002

Field Sample ID: MW-3

SAMPLE DATA

Lab Sample ID: 56637-21
Matrix: Aqueous
Percent Solid: N/A
Dilution Factor: 1.0
Collection Date: 05/26/06
Lab Receipt Date: 05/27/06
Extraction Date: 05/30/06
Analysis Date: 05/30/06

ANALYTICAL RESULTS SEMI-VOLATILE ORGANICS

ACID COMPOUND	Quantitation Limit µg/L	Result µg/L	ACID COMPOUND	Quantitation Limit µg/L	Result µg/L
2-Chlorophenol	5	U	Pentachlorophenol	10	U
4-Chloro-3-methylphenol	10	U	Phenol	5	U
2,4-Dichlorophenol	5	U	2,4,5-Trichlorophenol	5	U
2,4-Dimethylphenol	5	U	2,4,6-Trichlorophenol	5	U
2,4-dinitrophenol	5	U	Benzoic Acid	10	U
4,6-Dinitro-2-methylphenol	5	U	2-Methylphenol	5	U
2-Nitrophenol	5	U	3+4-Methylphenol	5	U
2,6-Dichlorophenol	5	U	Benzyl Alcohol	5	U
4-Nitrophenol	5	U	2,3,4,6-Tetrachlorophenol	5	U

Acid Surrogate Standard Recovery

2-Fluorophenol	44 %	d5-Phenol	33 %	2,4,6-Tribromophenol	77 %

BASE NEUTRAL COMPOUND	Quantitation Limit µg/L	Result µg/L	BASE NEUTRAL COMPOUND	Quantitation Limit µg/L	Result µg/L
1,2-Dichlorobenzene	2	U	Hexachlorobenzene	2	U
1,3-Dichlorobenzene	2	U	* Benzidine	20	U
1,4-Dichlorobenzene	2	U	3,3'-Dichlorobenzidine	20	U
2,4-Dinitrotoluene	2	U	Azobenzene	2	U
2,6-Dinitrotoluene	2	U	Bis(2-chloroethoxy)methane	2	U
Nitrobenzene	2	U	bis(2-chloroethyl) ether	2	U
Hexachlorobutadiene	2	U	bis(2-chloroisopropyl)ether	2	U
Dimethyl Phthalate	2	U	4-bromophenyl phenyl ether	2	U
Di-n-butyl phthalate	2	U	Butyl benzyl phthalate	2	U
di-n-octyl-phthalate	2	U	4-Chlorophenyl phenyl ether	2	U
Bis (2-ethylhexyl) phthalate	2	2	Diethyl Phthalate	2	U
1,2,4-Trichlorobenzene	2	U	Hexachlorocyclopentadiene	2	U

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis was conducted according to: Test Methods for Evaluating Solid Waste, SW-846 Method 8270C.

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8270/625 layout

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May 31, 2006

SAMPLE DATA

CLIENT SAMPLE ID

Project Name:	GP Old Town Mill
Project Number:	218736.002
Field Sample ID:	MW-3

Lab Sample ID:	56637-21
Matrix:	Aqueous
Percent Solid:	N/A
Dilution Factor:	1.0
Collection Date:	05/26/06
Lab Receipt Date:	05/27/06
Extraction Date:	05/30/06
Analysis Date:	05/30/06

ANALYTICAL RESULTS SEMI-VOLATILE ORGANICS

BASE NEUTRAL COMPOUND	Quantitation Limit µg/L	Result µg/L	BASE NEUTRAL COMPOUND	Quantitation Limit µg/L	Result µg/L
Acenaphthene	2	U	N-nitrosodimethylamine	2	U
Acenaphthylene	2	U	N-nitroso-di-n-propylamine	2	U
Anthracene	2	U	n-nitrosodiphenylamine	2	U
Benzo[a]anthracene	2	U	Pyridine	2	U
Benzo[a] pyrene	2	U	2-Methylnaphthalene	2	U
Benzo[b] fluoranthene	2	U	2-Chloronaphthalene	2	U
Benzo[k] fluoranthene	2	U	Naphthalene	2	U
Benzo(g,h,i) perylene	2	U	Phenanthrene	2	U
Chrysene	2	U	Dibenzofuran	2	U
Dibenz [a,h] anthracene	2	U	Aniline	2	U
Fluoranthene	2	U	4-Chloroaniline	2	U
Fluorene	2	U	2-Nitroaniline	2	U
Indeno [1,2,3-cd] pyrene	2	U	3-Nitroaniline	2	U
Pyrene	2	U	4-Nitroaniline	2	U
Hexachloroethane	2	U	Carbazole	2	U
Isophorone	2	U			

Base Neutral Surrogate Standard Recovery

2-Fluorobiphenyl	72	%	d5-nitrobenzene	71	%	d14-p-terphenyl	31 * %

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis was conducted according to: Test Methods for Evaluating Solid Waste, SW-846 Method 8270C.

COMMENTS: *Due to the reactive nature of this compound, the Benzidine quantitation limit is estimated.
*Surrogate recovery outside laboratory acceptance criteria. Sample was reanalyzed with similar results.

8270/625 layout

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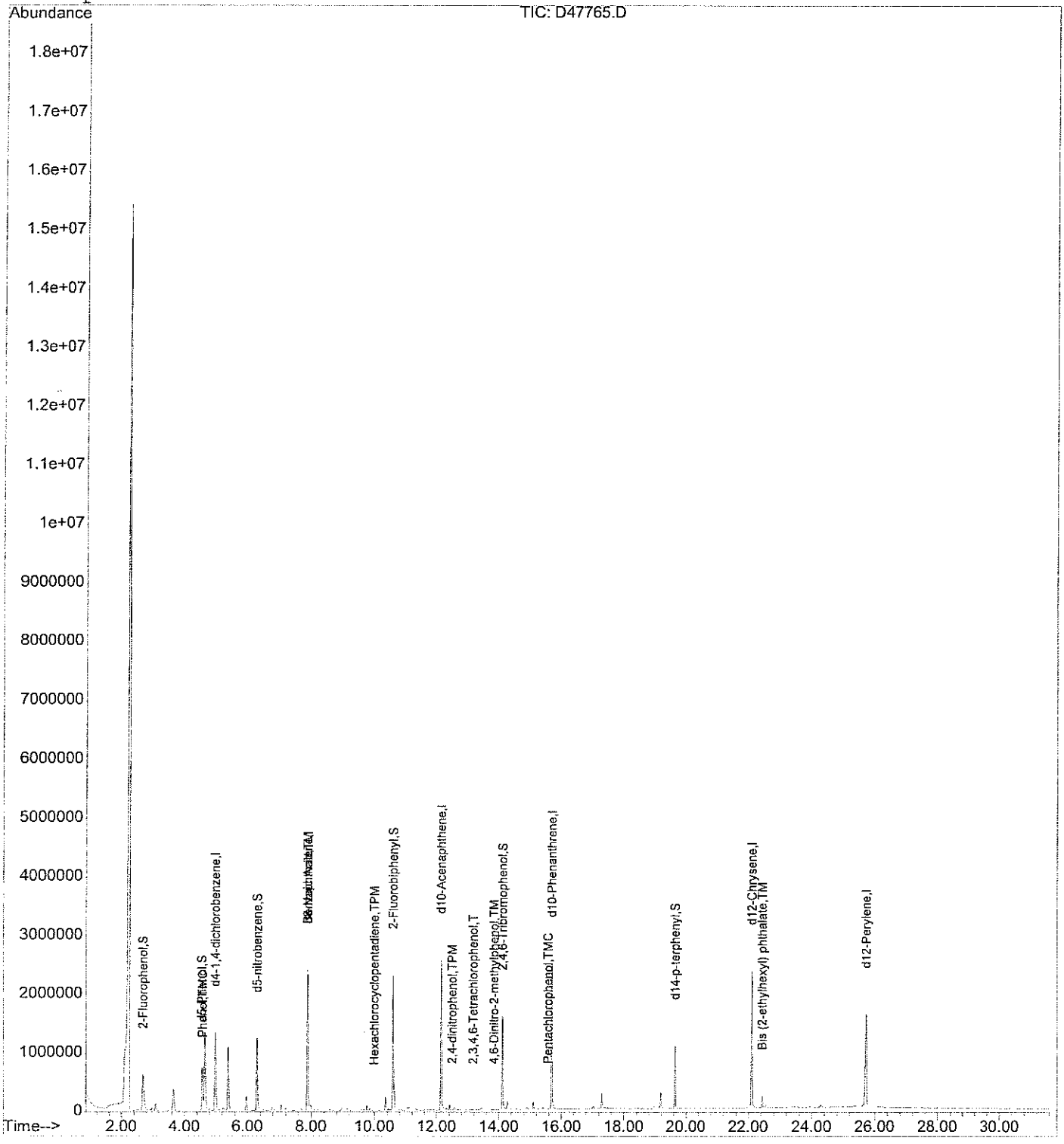
Quantitation Report

Data File : D:\HPCHEM\1\DATA\053006-D\D47765.D Vial: 11
Acq On : 30 May 2006 1:50 pm Operator:
Sample : 56637-21 Inst : D Instrum
Misc : Multiplr: 1.00
MS Integration Params: RTEINT.P
Quant Time: May 30 14:12 2006 Quant Results File: S05246.RES

Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Tue May 30 09:52:41 2006
Response via : Initial Calibration



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Portland ME 04102

May 31, 2006

CLIENT SAMPLE ID

Project Name: GP Old Town Mill

Project Number: 218736.002

Field Sample ID: Pile 1

SAMPLE DATA

Lab Sample ID: 56637-22
Matrix: Solid
Percent Solid: 90
Dilution Factor: 1.1
Collection Date: 05/26/06
Lab Receipt Date: 05/27/06
Extraction Date: 05/30/06
Analysis Date: 05/31/06

ANALYTICAL RESULTS SEMI-VOLATILE ORGANICS

ACID COMPOUND	Quantitation Limit µg/kg	Result µg/kg	ACID COMPOUND	Quantitation Limit µg/kg	Result µg/kg
2-Chlorophenol	270	U	Pentachlorophenol	270	U
4-Chloro-3-methylphenol	270	U	Phenol	270	U
2,4-Dichlorophenol	270	U	2,4,5-Trichlorophenol	270	U
2,4-Dimethylphenol	270	U	2,4,6-Trichlorophenol	270	U
2,4-dinitrophenol	270	U	Benzoic Acid	270	U
4,6-Dinitro-2-methylphenol	270	U	2-Methylphenol	270	U
2-Nitrophenol	270	U	3+4-Methylphenol	270	U
2,6-Dichlorophenol	270	U	Benzyl Alcohol	270	U
4-Nitrophenol	270	U	2,3,4,6-Tetrachlorophenol	270	U

Acid Surrogate Standard Recovery

2-Fluorophenol	65 %	d5-Phenol	80 %	2,4,6-Tribromophenol	87 %

BASE NEUTRAL COMPOUND	Quantitation Limit µg/kg	Result µg/kg	BASE NEUTRAL COMPOUND	Quantitation Limit µg/kg	Result µg/kg
1,2-Dichlorobenzene	270	U	Hexachlorobenzene	270	U
1,3-Dichlorobenzene	270	U	* Benzidine	270	U
1,4-Dichlorobenzene	270	U	3,3'-Dichlorobenzidine	270	U
2,4-Dinitrotoluene	270	U	Azobenzene	270	U
2,6-Dinitrotoluene	270	U	Bis(2-chloroethoxy)methane	270	U
Nitrobenzene	270	U	bis(2-chloroethyl) ether	270	U
Hexachlorobutadiene	270	U	bis(2-chloroisopropyl)ether	270	U
Dimethyl Phthalate	270	U	4-bromophenyl phenyl ether	270	U
Di-n-butyl phthalate	270	U	Butyl benzyl phthalate	270	U
di-n-octyl-phthalate	270	U	4-Chlorophenyl phenyl ether	270	U
Bis (2-ethylhexyl) phthalate	270	U	Diethyl Phthalate	270	U
1,2,4-Trichlorobenzene	270	U	Hexachlorocyclopentadiene	270	U

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis was conducted according to: Test Methods for Evaluating Solid Waste, SW-846 Method 8270C.

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8270/625 layout

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June 1, 2006

CLIENT SAMPLE ID

Project Name: GP Old Town Mill

Project Number: 218736.002

Field Sample ID: Pile 1

SAMPLE DATA

Lab Sample ID: 56637-22
Matrix: Solid
Percent Solid: 90
Dilution Factor: 1.1
Collection Date: 05/26/06
Lab Receipt Date: 05/27/06
Extraction Date: 05/30/06
Analysis Date: 05/31/06

ANALYTICAL RESULTS SEMI-VOLATILE ORGANICS

BASE NEUTRAL COMPOUND	Quantitation Limit µg/kg	Result µg/kg	BASE NEUTRAL COMPOUND	Quantitation Limit µg/kg	Result µg/kg
Acenaphthene	270	U	N-nitrosodimethylamine	270	U
Acenaphthylene	270	**216 J**	N-nitroso-di-n-propylamine	270	U
Anthracene	270	**160 J**	n-nitrosodiphenylamine	270	U
Benzo[a]anthracene	270	**567**	Pyridine	270	U
Benzo[a] pyrene	270	**666**	2-Methylnaphthalene	270	U
Benzo[b] fluoranthene	270	**585**	2-Chloronaphthalene	270	U
Benzo[k] fluoranthene	270	**590**	Naphthalene	270	U
Benzo(g,h,i) perylene	270	**140 J**	Phenanthrene	270	**691**
Chrysene	270	**720**	Dibenzofuran	270	U
Dibenz [a,h] anthracene	270	U	Aniline	270	U
Fluoranthene	270	**1080**	4-Chloroaniline	270	U
Fluorene	270	U	2-Nitroaniline	270	U
Indeno [1,2,3-cd] pyrene	270	**149 J**	3-Nitroaniline	270	U
Pyrene	270	**1730**	4-Nitroaniline	270	U
Hexachloroethane	270	U	Carbazole	270	U
Isophorone	270	U			

Base Neutral Surrogate Standard Recovery

2-Fluorobiphenyl	79	%	d5-nitrobenzene	74	%	d14-p-terphenyl	97 %

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis was conducted according to: Test Methods for Evaluating Solid Waste, SW-846 Method 8270C.

COMMENTS: *Due to the reactive nature of this compound, the Benzidine quantitation limit is estimated.
Results are expressed on a dry weight basis. Sample had low recovery of internal standard d12-Perylene. Sample was reanlayzed at a dilution with acceptable results.

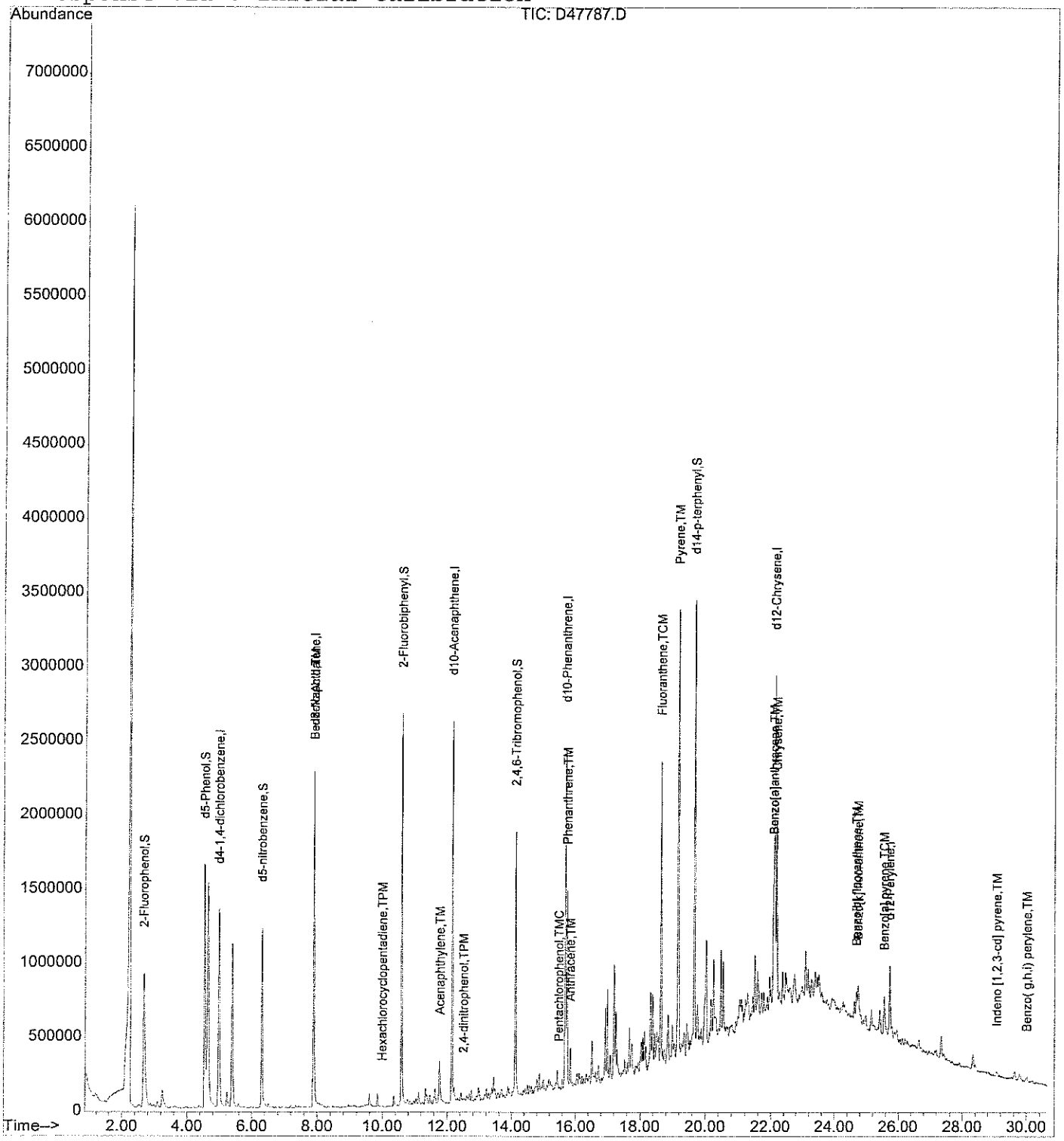
8270/625 layout

Authorized signature Melena Hall

Quantitation Report

Data File : D:\HPCHEM\1\DATA\053006-D\D47787.D Vial: 17
Acq On : 31 May 2006 5:49 am Operator:
Sample : 56637-22 Inst : D Instrum
Misc : SOIL Multiplr: 1.00
MS Integration Params: RTEINT.P
Quant Time: May 31 8:21 2006 Quant Results File: S05246.RES

Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Tue May 30 20:31:56 2006
Response via : Initial Calibration



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June 1, 2006

CLIENT SAMPLE ID

Project Name: GP Old Town Mill

Project Number: 218736.002

Field Sample ID: Pile 1

SAMPLE DATA

Lab Sample ID: 56637-22 DL
Matrix: Solid
Percent Solid: 90
Dilution Factor: 1.1
Collection Date: 05/26/06
Lab Receipt Date: 05/27/06
Extraction Date: 05/30/06
Analysis Date: 05/31/06

ANALYTICAL RESULTS SEMI-VOLATILE ORGANICS

ACID COMPOUND	Quantitation Limit µg/kg	Result µg/kg	ACID COMPOUND	Quantitation Limit µg/kg	Result µg/kg
2-Chlorophenol	270	U	Pentachlorophenol	270	U
4-Chloro-3-methylphenol	270	U	Phenol	270	U
2,4-Dichlorophenol	270	U	2,4,5-Trichlorophenol	270	U
2,4-Dimethylphenol	270	U	2,4,6-Trichlorophenol	270	U
2,4-dinitrophenol	270	U	Benzoic Acid	270	U
4,6-Dinitro-2-methylphenol	270	U	2-Methylphenol	270	U
2-Nitrophenol	270	U	3+4-Methylphenol	270	U
2,6-Dichlorophenol	270	U	Benzyl Alcohol	270	U
4-Nitrophenol	270	U	2,3,4,6-Tetrachlorophenol	270	U

Acid Surrogate Standard Recovery

2-Fluorophenol	69 %	d5-Phenol	86 %	2,4,6-Tribromophenol	101 %

BASE NEUTRAL COMPOUND	Quantitation Limit µg/kg	Result µg/kg	BASE NEUTRAL COMPOUND	Quantitation Limit µg/kg	Result µg/kg
1,2-Dichlorobenzene	270	U	Hexachlorobenzene	270	U
1,3-Dichlorobenzene	270	U	* Benzidine	270	U
1,4-Dichlorobenzene	270	U	3,3'-Dichlorobenzidine	270	U
2,4-Dinitrotoluene	270	U	Azobenzene	270	U
2,6-Dinitrotoluene	270	U	Bis(2-chloroethoxy)methane	270	U
Nitrobenzene	270	U	bis(2-chloroethyl) ether	270	U
Hexachlorobutadiene	270	U	bis(2-chloroisopropyl)ether	270	U
Dimethyl Phthalate	270	U	4-bromophenyl phenyl ether	270	U
Di-n-butyl phthalate	270	U	Butyl benzyl phthalate	270	U
di-n-octyl-phthalate	270	U	4-Chlorophenyl phenyl ether	270	U
Bis (2-ethylhexyl) phthalate	270	U	Diethyl Phthalate	270	U
1,2,4-Trichlorobenzene	270	U	Hexachlorocyclopentadiene	270	U

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis was conducted according to: Test Methods for Evaluating Solid Waste, SW-846 Method 8270C.

Authorized signature Melena Lalli

8270/625 layout

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June 1, 2006

CLIENT SAMPLE ID

Project Name: GP Old Town Mill

Project Number: 218736.002

Field Sample ID: Pile 1

SAMPLE DATA

Lab Sample ID:	56637-22	DL
Matrix:	Solid	
Percent Solid:	90	
Dilution Factor:	1.1	
Collection Date:	05/26/06	
Lab Receipt Date:	05/27/06	
Extraction Date:	05/30/06	
Analysis Date:	05/31/06	

ANALYTICAL RESULTS SEMI-VOLATILE ORGANICS

BASE NEUTRAL COMPOUND	Quantitation Limit µg/kg	Result µg/kg	BASE NEUTRAL COMPOUND	Quantitation Limit µg/kg	Result µg/kg
Acenaphthene	270	U	N-nitrosodimethylamine	270	U
Acenaphthylene	270	U	N-nitroso-di-n-propylamine	270	U
Anthracene	270	U	n-nitrosodiphenylamine	270	U
Benzo[a]anthracene	270	**278**	Pyridine	270	U
Benzo[a] pyrene	270	**364**	2-Methylnaphthalene	270	U
Benzo[b] fluoranthene	270	**286**	2-Chloronaphthalene	270	U
Benzo[k] fluoranthene	270	**289**	Naphthalene	270	U
Benzo(g,h,i) perylene	270	U	Phenanthrene	270	**377**
Chrysene	270	**349**	Dibenzofuran	270	U
Dibenz [a,h] anthracene	270	U	Aniline	270	U
Fluoranthene	270	**572**	4-Chloroaniline	270	U
Fluorene	270	U	2-Nitroaniline	270	U
Indeno [1,2,3-cd] pyrene	270	U	3-Nitroaniline	270	U
Pyrene	270	**738**	4-Nitroaniline	270	U
Hexachloroethane	270	U	Carbazole	270	U
Isophorone	270	U			

Base Neutral Surrogate Standard Recovery

2-Fluorobiphenyl	87	%	d5-nitrobenzene	81	%	d14-p-terphenyl	79	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis was conducted according to: Test Methods for Evaluating Solid Waste, SW-846 Method 8270C.

COMMENTS: *Due to the reactive nature of this compound, the Benzidine quantitation limit is estimated.
Results are expressed on a dry weight basis.

8270/625 layout

Authorized signature

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May 31, 2006

CLIENT SAMPLE ID

Project Name: GP Old Town Mill

Project Number: 218736.002

Field Sample ID: Pile 2

SAMPLE DATA

Lab Sample ID: 56637-23
Matrix: Solid
Percent Solid: 97
Dilution Factor: 1.0
Collection Date: 05/26/06
Lab Receipt Date: 05/27/06
Extraction Date: 05/30/06
Analysis Date: 05/31/06

ANALYTICAL RESULTS SEMI-VOLATILE ORGANICS

ACID COMPOUND	Quantitation Limit µg/kg	Result µg/kg	ACID COMPOUND	Quantitation Limit µg/kg	Result µg/kg
2-Chlorophenol	250	U	Pentachlorophenol	250	U
4-Chloro-3-methylphenol	250	U	Phenol	250	U
2,4-Dichlorophenol	250	U	2,4,5-Trichlorophenol	250	U
2,4-Dimethylphenol	250	U	2,4,6-Trichlorophenol	250	U
2,4-dinitrophenol	250	U	Benzoic Acid	250	U
4,6-Dinitro-2-methylphenol	250	U	2-Methylphenol	250	U
2-Nitrophenol	250	U	3+4-Methylphenol	250	U
2,6-Dichlorophenol	250	U	Benzyl Alcohol	250	U
4-Nitrophenol	250	U	2,3,4,6-Tetrachlorophenol	250	U

Acid Surrogate Standard Recovery

2-Fluorophenol	69 %	d5-Phenol	80 %	2,4,6-Tribromophenol	90 %

BASE NEUTRAL COMPOUND	Quantitation Limit µg/kg	Result µg/kg	BASE NEUTRAL COMPOUND	Quantitation Limit µg/kg	Result µg/kg
1,2-Dichlorobenzene	250	U	Hexachlorobenzene	250	U
1,3-Dichlorobenzene	250	U	* Benzidine	250	U
1,4-Dichlorobenzene	250	U	3,3'-Dichlorobenzidine	250	U
2,4-Dinitrotoluene	250	U	Azobenzene	250	U
2,6-Dinitrotoluene	250	U	Bis(2-chloroethoxy)methane	250	U
Nitrobenzene	250	U	bis(2-chloroethyl) ether	250	U
Hexachlorobutadiene	250	U	bis(2-chloroisopropyl)ether	250	U
Dimethyl Phthalate	250	U	4-bromophenyl phenyl ether	250	U
Di-n-butyl phthalate	250	U	Butyl benzyl phthalate	250	U
di-n-octyl-phthalate	250	U	4-Chlorophenyl phenyl ether	250	U
Bis (2-ethylhexyl) phthalate	250	U	Diethyl Phthalate	250	U
1,2,4-Trichlorobenzene	250	U	Hexachlorocyclopentadiene	250	U

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis was conducted according to: Test Methods for Evaluating Solid Waste, SW-846 Method 8270C.

Authorized signature

8270/625 layout

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May 31, 2006

CLIENT SAMPLE ID

Project Name: GP Old Town Mill

Project Number: 218736.002

Field Sample ID: Pile 2

SAMPLE DATA

Lab Sample ID: 56637-23
Matrix: Solid
Percent Solid: 97
Dilution Factor: 1.0
Collection Date: 05/26/06
Lab Receipt Date: 05/27/06
Extraction Date: 05/30/06
Analysis Date: 05/31/06

ANALYTICAL RESULTS SEMI-VOLATILE ORGANICS

BASE NEUTRAL COMPOUND	Quantitation Limit µg/kg	Result µg/kg	BASE NEUTRAL COMPOUND	Quantitation Limit µg/kg	Result µg/kg
Acenaphthene	250	U	N-nitrosodimethylamine	250	U
Acenaphthylene	250	U	N-nitroso-di-n-propylamine	250	U
Anthracene	250	U	n-nitrosodiphenylamine	250	U
Benzo[a]anthracene	250	**179 J**	Pyridine	250	U
Benzo[a] pyrene	250	**199 J**	2-Methylnaphthalene	250	U
Benzo[b] fluoranthene	250	**342**	2-Chloronaphthalene	250	U
Benzo[k] fluoranthene	250	**248 J**	Naphthalene	250	U
Benzo(g,h,i) perylene	250	U	Phenanthrene	250	U
Chrysene	250	**222 J**	Dibenzofuran	250	U
Dibenz [a,h] anthracene	250	U	Aniline	250	U
Fluoranthene	250	**276**	4-Chloroaniline	250	U
Fluorene	250	U	2-Nitroaniline	250	U
Indeno [1,2,3-cd] pyrene	250	U	3-Nitroaniline	250	U
Pyrene	250	**322**	4-Nitroaniline	250	U
Hexachloroethane	250	U	Carbazole	250	U
Isophorone	250	U			

Base Neutral Surrogate Standard Recovery

2-Fluorobiphenyl	83 %	d5-nitrobenzene	72 %	d14-p-terphenyl	92 %

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis was conducted according to: Test Methods for Evaluating Solid Waste, SW-846 Method 8270C.

COMMENTS: *Due to the reactive nature of this compound, the Benzidine quantitation limit is estimated.
Results are expressed on a dry weight basis. Sample had low recovery of internal standard d12-Perylene. Sample is being reanlayzed to confirm results.

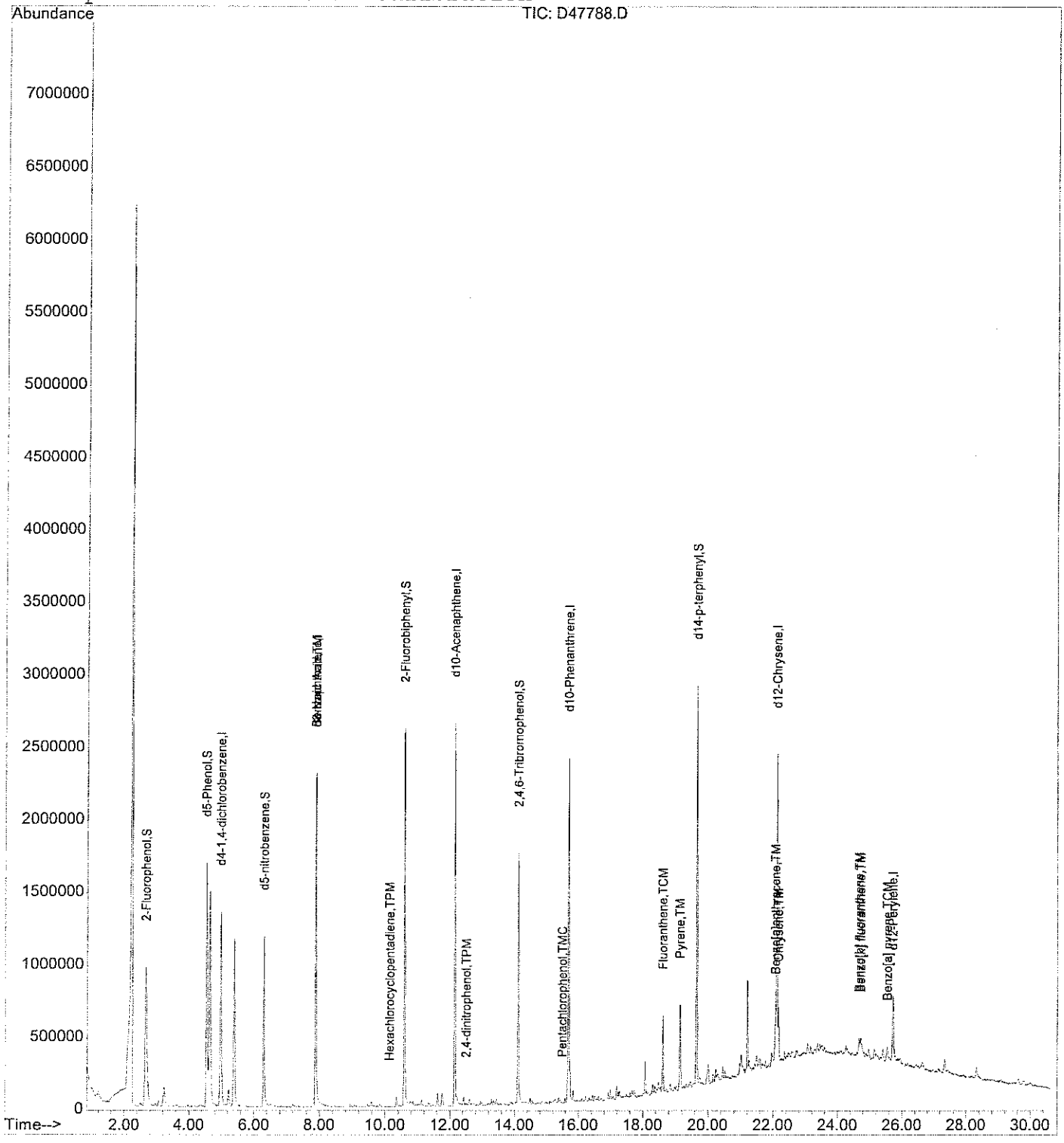
8270/625 layout

Authorized signature Melissa Tull

Quantitation Report

Data File : D:\HPCHEM\1\DATA\053006-D\D47788.D Vial: 18
Acq On : 31 May 2006 6:22 am Operator:
Sample : 56637-23 Inst : D Instrum
Misc : SOIL Multiplr: 1.00
MS Integration Params: RTEINT.P
Quant Time: May 31 8:21 2006 Quant Results File: S05246.RES

Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Tue May 30 20:31:56 2006
Response via : Initial Calibration



analytics environmental laboratory LLC

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800-929-9906

Mr. Karl Kasper
Woodard & Curran
41 Hutchins Drive
Portland ME 04102

June 1, 2006

CLIENT SAMPLE ID

Project Name: GP Old Town Mill

Project Number: 218736.002

Field Sample ID: Pile 2

SAMPLE DATA

Lab Sample ID: 56637-23 DL
Matrix: Solid
Percent Solid: 97
Dilution Factor: 2.0
Collection Date: 05/26/06
Lab Receipt Date: 05/27/06
Extraction Date: 05/30/06
Analysis Date: 05/31/06

ANALYTICAL RESULTS SEMI-VOLATILE ORGANICS

ACID COMPOUND	Quantitation Limit µg/kg	Result µg/kg	ACID COMPOUND	Quantitation Limit µg/kg	Result µg/kg
2-Chlorophenol	510	U	Pentachlorophenol	510	U
4-Chloro-3-methylphenol	510	U	Phenol	510	U
2,4-Dichlorophenol	510	U	2,4,5-Trichlorophenol	510	U
2,4-Dimethylphenol	510	U	2,4,6-Trichlorophenol	510	U
2,4-dinitrophenol	510	U	Benzoic Acid	510	U
4,6-Dinitro-2-methylphenol	510	U	2-Methylphenol	510	U
2-Nitrophenol	510	U	3+4-Methylphenol	510	U
2,6-Dichlorophenol	510	U	Benzyl Alcohol	510	U
4-Nitrophenol	510	U	2,3,4,6-Tetrachlorophenol	510	U

Acid Surrogate Standard Recovery

2-Fluorophenol	72 %	d5-Phenol	84 %	2,4,6-Tribromophenol	99 %

BASE NEUTRAL COMPOUND	Quantitation Limit µg/kg	Result µg/kg	BASE NEUTRAL COMPOUND	Quantitation Limit µg/kg	Result µg/kg
1,2-Dichlorobenzene	510	U	Hexachlorobenzene	510	U
1,3-Dichlorobenzene	510	U	* Benzidine	510	U
1,4-Dichlorobenzene	510	U	3,3'-Dichlorobenzidine	510	U
2,4-Dinitrotoluene	510	U	Azobenzene	510	U
2,6-Dinitrotoluene	510	U	Bis(2-chloroethoxy)methane	510	U
Nitrobenzene	510	U	bis(2-chloroethyl) ether	510	U
Hexachlorobutadiene	510	U	bis(2-chloroisopropyl)ether	510	U
Dimethyl Phthalate	510	U	4-bromophenyl phenyl ether	510	U
Di-n-butyl phthalate	510	U	Butyl benzyl phthalate	510	U
di-n-octyl-phthalate	510	U	4-Chlorophenyl phenyl ether	510	U
Bis (2-ethylhexyl) phthalate	510	U	Diethyl Phthalate	510	U
1,2,4-Trichlorobenzene	510	U	Hexachlorocyclopentadiene	510	U

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis was conducted according to: Test Methods for Evaluating Solid Waste, SW-846 Method 8270C.

Authorized signature

8270/625 layout

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Mr. Karl Kasper
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41 Hutchins Drive
Portland ME 04102

June 1, 2006

CLIENT SAMPLE ID

Project Name: GP Old Town Mill

Project Number: 218736.002

Field Sample ID: Pile 2

SAMPLE DATA

Lab Sample ID:	56637-23	DL
Matrix:	Solid	
Percent Solid:	97	
Dilution Factor:	2.0	
Collection Date:	05/26/06	
Lab Receipt Date:	05/27/06	
Extraction Date:	05/30/06	
Analysis Date:	05/31/06	

ANALYTICAL RESULTS SEMI-VOLATILE ORGANICS

BASE NEUTRAL COMPOUND	Quantitation Limit µg/kg	Result µg/kg	BASE NEUTRAL COMPOUND	Quantitation Limit µg/kg	Result µg/kg
Acenaphthene	510	U	N-nitrosodimethylamine	510	U
Acenaphthylene	510	U	N-nitroso-di-n-propylamine	510	U
Anthracene	510	U	n-nitrosodiphenylamine	510	U
Benzo[a]anthracene	510	U	Pyridine	510	U
Benzo[a] pyrene	510	U	2-Methylnaphthalene	510	U
Benzo[b] fluoranthene	510	U	2-Chloronaphthalene	510	U
Benzo[k] fluoranthene	510	U	Naphthalene	510	U
Benzo(g,h,i) perylene	510	U	Phenanthrene	510	U
Chrysene	510	U	Dibenzofuran	510	U
Dibenz [a,h] anthracene	510	U	Aniline	510	U
Fluoranthene	510	**300 J**	4-Chloroaniline	510	U
Fluorene	510	U	2-Nitroaniline	510	U
Indeno [1,2,3-cd] pyrene	510	U	3-Nitroaniline	510	U
Pyrene	510	**314 J**	4-Nitroaniline	510	U
Hexachloroethane	510	U	Carbazole	510	U
Isophorone	510	U			

Base Neutral Surrogate Standard Recovery

2-Fluorobiphenyl	84	%	d5-nitrobenzene	76	%	d14-p-terphenyl	86 %

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis was conducted according to: Test Methods for Evaluating Solid Waste, SW-846 Method 8270C.

COMMENTS: *Due to the reactive nature of this compound, the Benzidine quantitation limit is estimated.
Results are expressed on a dry weight basis.

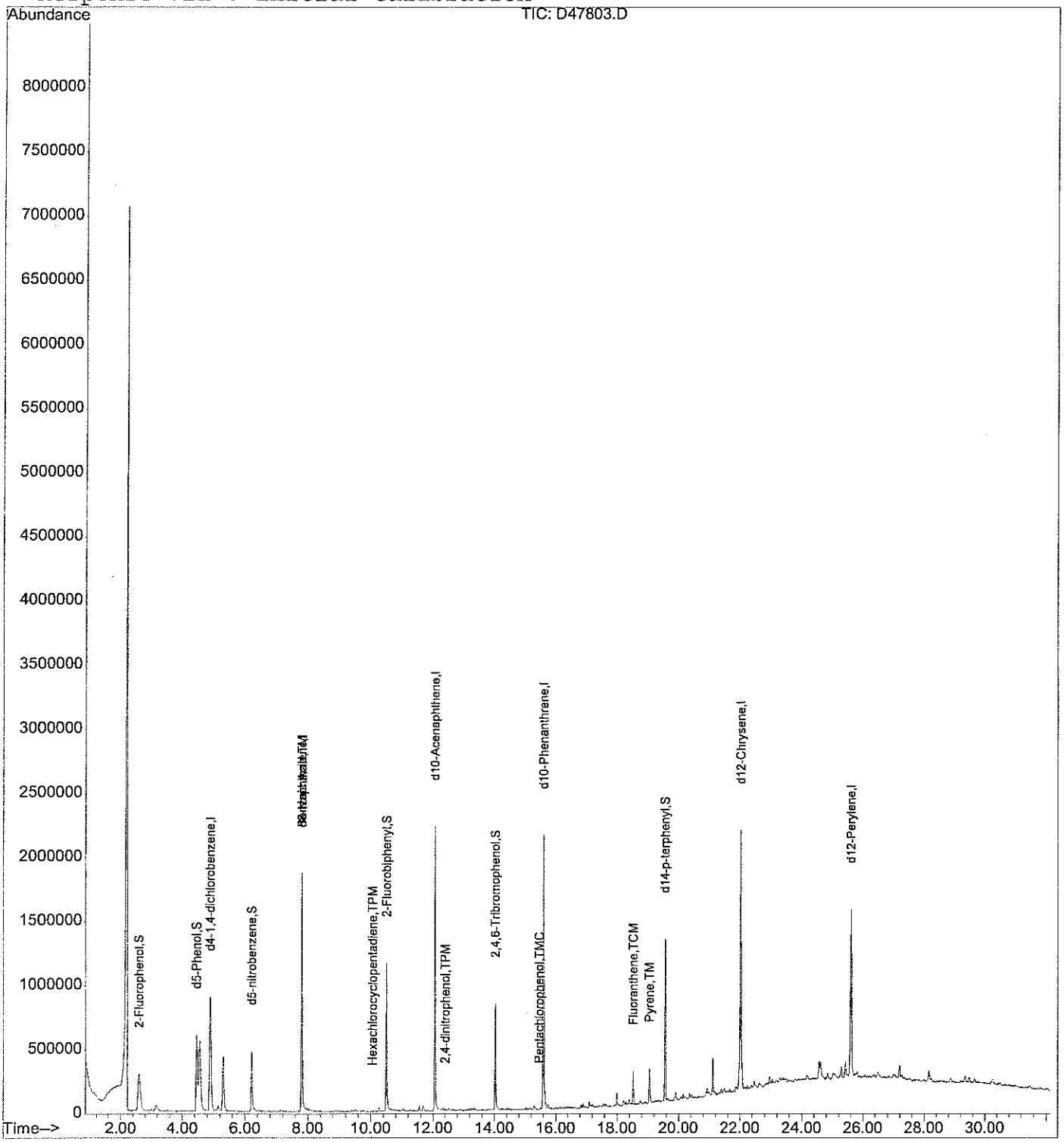
8270/625 layout

Authorized signature

Quantitation Report

Data File : D:\HPCHEM\1\DATA\053106-D\D47803.D Vial: 10
Acq On : 31 May 2006 4:10 pm Operator:
Sample : 56637-23,,1:2 Inst : D Instrum
Misc : SOIL Multiplr: 1.00
MS Integration Params: RTEINT.P
Quant Time: May 31 19:58 2006 Quant Results File: S05246.RES

Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Wed May 31 19:58:02 2006
Response via : Initial Calibration



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SEMI-VOLATILE
QC FORMS

AnalyticsLLC:AEL Documents LLC:Pkg Dividers:SVOCQC.doc

analytics environmental laboratory LLC

SEMI-VOLATILE ORGANIC SOIL
SYSTEM MONITORING COMPOUNDS
SUMMARY

Project #: 218736.002 Quote #: NA SDG#: 56637

DFTPP File ID: D47722T.D Site ID: GP Old Town Mill Location: ME

Instrument ID: D DFTPP Injection Date: 5/25/06

GC Column: RTX-5MS ID: 0.25(mm) DFTPP Injection Time: 0305

	SAMPLE ID	SMC1 #	SMC 2 #	SMC 3 #	SMC 4 #	SMC 5 #	SMC 6 #	TOT OUT
01	S6852 PT1 2/4	NA	NA	NA	NA	NA	NA	0
02	S6851 PT2A 4/8	NA	NA	NA	NA	NA	NA	0
03	S6850 PT2 10/20	NA	NA	NA	NA	NA	NA	0
04	S6849 PT3 20/40	44	44	81	81	33 *	79	0
05	S6848 PT4A 40/80	88	90	159 *	164 *	71	160 *	0
06	S6847 PT4 50/100	NA	NA	NA	NA	NA	NA	0
07	S6846 PT5 100/200	NA	NA	NA	NA	NA	NA	0
08	S6857 ABN CC 20/40	41	30 *	83	82	41	77	0
09	S6858 BENZIDINES + ANIL	NA	NA	NA	NA	NA	NA	0
10								
11								
12								
13								
14								
15								
16								
17								
18								
19								
20								
21								
22								
23								
24								
25								
26								

	QC LIMITS
SMC1 = 2-Fluorophe	29-97 %
SMC2 = d5-Phenol	40-120 %
SMC3 = d5-Nitrobenzene	20-101%
SMC4 = 2-Fluorobiphenyl	34-102 %
SMC5 = 2,4,6-Tribromophenol	36-141 %
SMC6 = d14-p-Terphenyl	38-108%

Column to be used to flag recovery values

* Values outside of protocol required QC limits

D System Monitoring Compound diluted out

analytics environmental laboratory LLC

SEMI-VOLATILE ORGANIC AQUEOUS SYSTEM MONITORING COMPOUNDS SUMMARY

Project #: 218736.002 Quote #: NA SDG#: 56637

DFTPP File ID: D47755T.D Site ID: GP Old Town Mill Location: ME

Instrument ID: D DFTPP Injection Date: 5/30/06

GC Column: RTX-5MS ID: 0.25(mm) DFTPP Injection Time: 0639

	SAMPLE ID	SMC1 #	SMC 2 #	SMC 3 #	SMC 4 #	SMC 5 #	SMC 6 #	TOT OUT
01	S6849 PT3 20/40	40	40	83	79	42 *	78	1
02	S6760 B+A 20	NA	NA	NA	NA	NA	NA	0
03	B05306AW	51	37	81	75	85	84	0
04	L05306AW	47	34	79	77	82	83	0
05	LD05306AW	43	31	75	75	76	78	0
06	56637-1	38	31	77	75	62 *	48	1
07	56637-2	2 *	0 *	76	74	8 *	55	3
08	56637-3	40	29	81	83	82	73	0
09	56637-20	2 *	1 *	70	72	5 *	73	3
10	56637-21	44	33	71	72	77	31 *	1
11	56637-1, RE-INJECT	38	30	76	76	60 *	48	1
12	56637-2, RE-INJECT	2 *	0 *	76	76	7 *	54	3
13	56637-20, RE-INJE(	2 *	0 *	71	73	5 *	74	3
14	56637-21, RE-INJE(	44	32	74	72	75	31 *	1
15	LD05306AW, RE-INJECT	43	31	75	75	75	80	0
16								
17								
18								
19								
20								
21								
22								
23								
24								
25								
26								

	QC LIMITS
SMC1 = 2-Fluorophenol	28-110%
SMC2 = d5-Phenol	22-110%
SMC3 = d5-Nitrobenzene	45-107%
SMC4 = 2-Fluorobiphenyl	44-102%
SMC5 = 2,4,6-Tribromophenol	63-127%
SMC6 = d14-p-Terphenyl	33-132%

Column to be used to flag recovery values

* Values outside of protocol required QC limits

D System Monitoring Compound diluted out

analytics environmental laboratory LLC

SEMI-VOLATILE ORGANIC SOIL SYSTEM MONITORING COMPOUNDS SUMMARY

Project #:	218736.002	Quote #:	NA	SDG#:	56637
DFTPP File ID:	D47771T.D	Site ID:	GP Old Town Mill	Location:	ME
Instrument ID:	D			DFTPP Injection Date:	5/30/06
GC Column: RTX-5MS		ID: 0.25(mm)		DFTPP Injection Time:	2043

	SAMPLE ID	SMC1 #	SMC 2#	SMC 3 #	SMC 4 #	SMC 5 #	SMC 6 #	TOT OUT
01	S6849 PT3 20/40	52	52	82	79	51	82	0
02	S6760 B+A 20	NA	NA	NA	NA	NA	NA	0
03	B05306AS	68	77	72	77	77	83	0
04	L05306AS	65	74	73	77	85	84	0
05	LD05306AS	67	76	75	79	85	83	0
06	ZZZZZ	70	80	76	77	84	85	0
07	ZZZZZ	67	77	74	76	84	85	0
08	56637-4	68	80	74	77	90	93	0
09	56637-9	61	72	66	72	83	85	0
10	56637-10	56	67	59	68	87	88	0
11	56637-12	67	77	72	74	80	88	0
12	56637-18	79	89	87	79	94	91	0
13	56637-11	62	75	64	75	93	85	0
14	56637-14	65	77	71	77	90	77	0
15	56637-19	73	84	77	83	95	70	0
16	56637-22	65	80	74	79	87	97	0
17	56637-23	69	80	72	83	90	92	0
18	56637-15,,1:2	74	84	75	79	96	220 *	1
19	56637-16,,1:5	66	74	68	69	79	185 *	1
20	56637-17,,1:5	57	61	59	67	69	66	0
21								
22								
23								
24								
25								
26								

	QC LIMITS
SMC1 = 2-Fluorophenol	29-97 %
SMC2 = d5-Phenol	40-120 %
SMC3 = d5-Nitrobenzene	20-101%
SMC4 = 2-Fluorobiphenyl	34-102 %
SMC5 = 2,4,6-Tribromophenol	36-141 %
SMC6 = d14-p-Terphenyl	38-108%

\# Column to be used to flag recovery values

* Values outside of protocol required QC limits

D System Monitoring Compound diluted out

analytics environmental laboratory LLC

SEMI-VOLATILE ORGANIC SOIL
SYSTEM MONITORING COMPOUNDS
SUMMARY

Project #: 218736.002 Quote #: NA SDG#: 56637

DFTPP File ID: D47794T.D Site ID: GP Old Town Mill Location: ME

Instrument ID: D DFTPP Injection Date: 05/31/06

GC Column: RTX-5MS ID: 0.25(mm) DFTPP Injection Time: 10:18

	SAMPLE ID	SMC1 #	SMC 2 #	SMC 3 #	SMC 4 #	SMC 5 #	SMC 6 #	TOT OUT
01	S6849 PT3 20/40	53	53	82	81	56	84	0
02	S6740 B+A 20	NA	NA	NA	NA	NA	NA	0
03	B05306AS	67	78	72	77	85	86	0
04	56637-15,,1:100	D	D	D	D	D	D	0
05	56637-16,,1:50	D	D	D	D	D	D	0
06	56637-17,,1:100	D	D	D	D	D	D	0
07	56637-14,DL,1:2	68	81	74	82	92	66	0
08	56637-22,DL,1:2	69	86	81	87	101	79	0
09	56637-23,DL,1:2	72	84	76	84	99	86	0
10	56637-17,,1:20	D	D	D	D	D	D	0
11	56637-23,MS,,1:2	56	92	82	90	102	88	0
12	56637-23,MSD,,1:2	72	82	78	84	98	86	0
13	ZZZZZ	57	71	58	38	86	72	0
14	ZZZZZ	54	70	54	48	82	56	0
15	ZZZZZ	38	44	34	32 *	56	54	1
16	ZZZZZ	52	67	50	44	81	59	0
17								
18								
19								
20								
21								
22								
23								
24								
25								
26								

	QC LIMITS
SMC1 = 2-Fluorophenol	29-97 %
SMC2 = d5-Phenol	40-120 %
SMC3 = d5-Nitrobenzene	20-101%
SMC4 = 2-Fluorobiphenyl	34-102 %
SMC5 = 2,4,6-Tribromophenol	36-141 %
SMC6 = d14-p-Terphenyl	38-108%

Column to be used to flag recovery values

* Values outside of protocol required QC limits

D System Monitoring Compound diluted out

analytics environmental laboratory LLC

SEMI-VOLATILE ORGANIC AQUEOUS
LABORATORY CONTROL SPIKE/ DUPLICATE
PERCENT RECOVERY

Project #: 218736.002 SDG# 56637

DFTPP File ID: D47755T.D Site ID: Location: ME

Instrument ID: D DFTPP Injection Date: 5/30/06

GC Column: RTX-5MS ID: 0.25(mm) DFTPP Injection Time: 0639

Laboratory Control Spike/Duplicate Sample ID: 05306AW/LD05306AW

COMPOUND	SPIKE ADDED (ug/L)	SAMPLE AMT (ug/L)	MS CONC. (ug/L)	MS REC %	#	SPIKE ADDED (ug/L)	MSD CONC. (ug/L)	MSD REC %	#	RPD	#	QC LIMITS RPD	
Pyridine	40	U	18	46		40	3	7	*	147	*	25	27-56
N-nitrosodimethylamine	40	U	20	49		40	18	46		7		25	25-70
Aniline	40	U	29	73		40	7	18	*	121	*	25	38-84
Phenol	40	U	15	38		40	14	35		9		42	25-54
2-Chlorophenol	40	U	28	69		40	26	65		6		40	40-95
bis(2-chloroethyl) eth	40	U	32	80		40	29	73		9		25	40-104
1,3-Dichlorobenzene	40	U	26	66		40	24	61		8		25	35-75
1,4-Dichlorobenzene	40	U	27	67		40	24	61		8		28	37-75
1,2-Dichlorobenzene	40	U	28	69		40	25	63		9		25	37-79
Benzyl Alcohol	40	U	30	76		40	28	70		8		25	48-99
2-Methylphenol	40	U	27	67		40	25	62		8		25	48-90
bis(2-chloroisopropyl)ether	40	U	30	75		40	28	69		8		25	47-97
3+4-Methylphenol	40	U	26	65		40	24	60		9		25	46-85
Hexachloroethane	40	U	26	66		40	24	61		8		25	36-76
N-nitroso-di-n-propylamine	40	U	35	87		40	33	81		6		38	46-108
Nitrobenzene	40	U	33	83		40	31	78		6		25	52-92
Isophorone	40	U	39	97		40	37	93		4		25	50-110
2-Nitrophenol	40	U	30	76		40	28	71		7		25	41-103
2,4-Dimethylphenol	40	U	28	70		40	26	64		8		25	51-84
Bis(2-chloroethoxy)methane	40	U	33	83		40	32	79		5		25	55-96
2,4-Dichlorophenol	40	U	32	79		40	30	74		7		25	50-105
Benzoic Acid	40	U	7	18		40	8	20		11		25	10-54
1,2,4-Trichlorobenzer	40	U	30	74		40	28	70		6		28	43-79
Naphthalene	40	U	31	77		40	29	72		7		25	46-88
4-Chloroaniline	40	U	33	82		40	11	27	*	100	*	25	56-98
Hexachlorobutadiene	40	U	28	71		40	26	64		9		25	36-71
4-Chloro-3-methylphenol	40	U	32	80		40	30	74		8		42	59-104
2-Methylnaphthalene	40	U	32	79		40	29	73		7		25	52-96
Hexachlorocyclopentadiene	40	U	31	79		40	29	74		7		25	32-84
2,4,6-Trichlorophenol	40	U	31	78		40	29	72		7		50	50-115
2,4,5-Trichlorophenol	40	U	30	75		40	29	74		2		47	50-110
2-Chloronaphthalene	40	U	32	79		40	30	75		6		25	52-93
2-Nitroaniline	40	U	36	89		40	34	85		5		25	63-115
2,6-Dinitrotoluene	40	U	37	93		40	36	90		3		25	66-111
Dimethyl Phthalate	40	U	17	43		40	17	41		4		25	25-72
Acenaphthylene	40	U	34	84		40	32	80		5		25	51-104
3-Nitroaniline	40	U	37	93		40	18	45	*	70	*	25	58-122
Acenaphthene	40	U	34	84		40	32	79		6		31	52-99
2,4-dinitrophenol	40	U	25	63		40	27	67		7		25	21-118
Dibenzofuran	40	U	33	84		40	31	78		7		25	59-102
4-Nitrophenol	40	U	14	34		40	12	31		8		50	10-69

analytics environmental laboratory LLC

SEMI-VOLATILE ORGANIC AQUEOUS
LABORATORY CONTROL SPIKE/ DUPLICATE
PERCENT RECOVERY

Project #: 218736.002 SDG# 56637

DFTPP File ID: D47755T.D Site ID: Location: ME

Instrument ID: D DFTPP Injection Date: 5/30/06

GC Column: RTX-5MS ID: 0.25(mm) DFTPP Injection Time: 0639

Laboratory Control Spike/Duplicate Sample ID:05306AW/LD05306AW

COMPOUND	SPIKE ADDED (ug/L)	SAMPLE AMT (ug/L)	MS CONC. (ug/L)	MS REC %	#	SPIKE ADDED (ug/L)	MSD CONC. (ug/L)	MSD REC %	#	RPD	#	QC LIMITS RPD	
2,4-Dinitrotoluene	40	U	39	97		40	37	93		5		38	63-116
Fluorene	40	U	35	87		40	33	84		4		25	55-105
Diethyl Phthalate	40	U	31	76		40	29	72		6		25	40-93
4-Chlorophenyl phenyl ether	40	U	36	89		40	34	84		6		25	54-101
4-Nitroaniline	40	U	38	94		40	30	76		21		25	61-120
4,6-Dinitro-2-methylphen	40	U	30	76		40	29	73		4		25	40-130
n-nitrosodiphenylamine	40	U	39	98		40	36	89		9		25	53-110
Azobenzene	40	U	33	83		40	31	78		6		25	61-112
4-bromophenyl phenyl ether	40	U	34	85		40	32	81		6		25	58-101
Hexachlorobenzene	40	U	36	91		40	35	88		3		25	61-103
Pentachlorophenol	40	U	23	58		40	25	62		7		50	40-108
Phenanthrene	40	U	36	89		40	34	85		5		25	60-107
Anthracene	40	U	37	93		40	35	88		6		25	62-110
Carbazole	80	U	71	89		80	66	83		8		25	62-114
Di-n-butyl phthalate	40	U	35	87		40	33	82		6		25	56-109
Fluoranthene	40	U	36	90		40	34	85		6		25	62-112
Benzidine	80	U	91	114		80	6	7	*	177	*	25	10-131
Pyrene	40	U	36	91		40	35	87		4		31	56-106
Butyl benzyl phthalate	40	U	31	78		40	29	73		6		25	46-103
Benzo[a]anthracene	40	U	37	94		40	36	89		5		25	60-110
Chrysene	40	U	38	95		40	35	88		8		25	61-109
3,3'-Dichlorobenzidine	80	U	78	98		80	4	5	*	181	*	25	57-110
Bis (2-ethylhexyl) phthala	40	U	37	93		40	35	87		7		25	53-123
di-n-octyl-phthalate	40	U	33	82		40	33	82		0		25	52-135
Benzo[b] fluoranthene	40	U	36	91		40	35	88		4		25	60-114
Benzo[k] fluoranthene	40	U	36	90		40	35	87		3		25	58-115
Benzo[a] pyrene	40	U	35	88		40	34	85		3		25	61-110
Indeno [1,2,3-cd] pyrene	40	U	37	93		40	36	90		3		25	54-125
Dibenz [a,h] anthrace	40	U	38	96		40	37	94		3		25	55-125
Benzo(g,h,i) perylene	40	U	37	93		40	37	92		2		25	50-125

Column to be used to flag recovery and RPD values with an asterisk

* Values outside of QC limits
RPD: 6 out of 71outside limits
Spike Recovery: 6 out of 142 outside limits

Comments: _____

analytics environmental laboratory LLC

SEMI-VOLATILE ORGANIC SOIL
LABORATORY CONTROL SPIKE/ DUPLICATE
PERCENT RECOVERY

Project #: 218736.002 SDG# 56637

DFTPP File ID: D47794T Site ID: GP Old Town Mill Location: ME

Instrument ID: D DFTPP Injection Date: 5/31/06

GC Column: RTX-5MS ID: 0.25(mm) DFTPP Injection Time: 1010

Laboratory Control Spike/Duplicate Sample ID: L05306AS/LD05306AS

COMPOUND	SPIKE ADDED (ug/kg)	SAMPLE AMT (ug/kg)	LCS CONC. (ug/kg)	LCS REC %	#	LCSD CONC. (ug/kg)	LCSD REC %	#	RPD	#	QC LIMITS RPD	
Pyridine	1333	U	628	47		649	49		3		25	13-77
N-nitrosodimethylamine	1333	U	844	63		806	60		5		25	20-98
Aniline	1333	U	834	63		908	68		8		25	60-120
Phenol	1333	U	933	70		911	68		2		35	51-96
2-Chlorophenol	1333	U	890	67		885	66		1		50	44-92
bis(2-chloroethyl) eth	1333	U	919	69		859	64		7		25	36-85
1,3-Dichlorobenzene	1333	U	860	65		843	63		2		25	27-90
1,4-Dichlorobenzene	1333	U	847	64		856	64		1		27	27-90
1,2-Dichlorobenzene	1333	U	867	65		859	64		1		25	30-89
Benzyl Alcohol	1333	U	978	73		1002	75		2		25	45-135
2-Methylphenol	1333	U	956	72		922	69		4		25	45-135
bis(2-chloroisopropyl)ether	1333	U	868	65		856	64		1		25	41-102
3+4-Methylpheno	1333	U	979	73		950	71		3		25	45-135
Hexachloroethane	1333	U	900	67		895	67		1		25	30-93
N-nitroso-di-n-propylamine	1333	U	996	75		985	74		1		38	46-100
Nitrobenzene	1333	U	955	72		944	71		1		25	41-93
Isophorone	1333	U	1129	85		1112	83		1		25	62-116
2-Nitrophenol	1333	U	955	72		961	72		1		25	43-95
2,4-Dimethylphenol	1333	U	894	67		879	66		2		25	45-79
Bis(2-chloroethoxy)methane	1333	U	963	72		931	70		3		25	44-100
2,4-Dichlorophenol	1333	U	1017	76		987	74		3		25	47-95
Benzoic Acid	1333	U	750	56		691	52		8		25	30-150
1,2,4-Trichlorobenzei	1333	U	944	71		931	70		1		23	37-92
Naphthalene	1333	U	916	69		907	68		1		25	42-99
4-Chloroaniline	1333	U	935	70		998	75		7		25	60-120
Hexachlorobutadiene	1333	U	943	71		929	70		1		25	35-94
4-Chloro-3-methylphenol	1333	U	1053	79		1034	78		2		33	53-100
2-Methylnaphthalene	1333	U	972	73		987	74		2		25	60-120
Hexachlorocyclopentadiene	1333	U	1097	82		1092	82		0		25	36-97
2,4,6-Trichlorophenol	1333	U	1015	76		1008	76		1		50	56-100
2,4,5-Trichlorophenol	1333	U	1023	77		997	75		3		47	45-135
2-Chloronaphthalene	1333	U	976	73		951	71		3		25	49-96
2-Nitroaniline	1333	U	1106	83		1129	85		2		25	60-120
2,6-Dinitrotoluene	1333	U	1119	84		1072	80		4		25	60-101
Dimethyl Phthala	1333	U	1094	82		1052	79		4		25	53-105
Acenaphthylene	1333	U	1012	76		986	74		3		25	54-97
3-Nitroaniline	1333	U	1142	86		1167	88		2		25	60-120
Acenaphthene	1333	U	1013	76		987	74		3		19	55-96
2,4-dinitrophenol	1333	U	920	69		928	70		1		25	26-85
Dibenzofuran	1333	U	1034	78		1049	79		1		25	60-120
4-Nitrophenol	1333	U	1086	81		975	73		11		50	57-113

analytics environmental laboratory LLC

SEMI-VOLATILE ORGANIC SOIL
LABORATORY CONTROL SPIKE/ DUPLICATE
PERCENT RECOVERY

Project #: 218736.002 SDG# 56637

DFTPP File ID: D47794T Site ID: GP Old Town Mill Location: ME

Instrument ID: D DFTPP Injection Date: 5/31/06

GC Column: RTX-5MS ID: 0.25(mm) DFTPP Injection Time: 1010

Laboratory Control Spike/Duplicate Sample ID: L05306AS/LD05306AS

COMPOUND	SPIKE ADDED (ug/kg)	SAMPLE AMT (ug/kg)	LCS CONC. (ug/kg)	LCS REC %	#	LCSD CONC. (ug/kg)	LCSD REC %	#	RPD	#	QC RPD	LIMITS
2,4-Dinitrotoluene	1333	U	1145	86		1112	83		3		47	66-107
Fluorene	1333	U	1046	78		1020	76		2		25	60-101
Diethyl Phthalate	1333	U	1148	86		1115	84		3		25	57-111
4-Chlorophenyl phenyl ether	1333	U	1081	81		1043	78		4		25	54-103
4-Nitroaniline	1333	U	1147	86		1149	86		0		25	60-120
4,6-Dinitro-2-methylphenol	1333	U	978	73		980	73		0		25	47-110
n-nitrosodiphenylamine	1333	U	1160	87		1136	85		2		25	60-121
Azobenzene	1333	U	1000	75		971	73		3		25	63-115
4-bromophenyl phenyl ether	1333	U	1007	76		996	75		1		25	55-109
Hexachlorobenzene	1333	U	1061	80		1023	77		4		25	60-104
Pentachlorophenol	1333	U	858	64		827	62		4		47	32-115
Phenanthrene	1333	U	1053	79		1033	77		2		25	59-103
Anthracene	1333	U	1063	80		1071	80		1		25	60-107
Carbazole	2667	U	2020	76		2029	76		0		25	67-113
Di-n-butyl phthalate	1333	U	1048	79		1067	80		2		25	67-119
Fluoranthene	1333	U	1014	76		1032	77		2		25	65-114
Benzidine	2667	U	1294	49		1788	67		32		25	29-187
Pyrene	1333	U	1165	87		1082	81		7		36	52-110
Butyl benzyl phthalate	1333	U	1153	86		1085	81		6		25	56-121
Benzo[a]anthrace	1333	U	1151	86		1096	82		5		25	59-110
Chrysene	1333	U	1138	85		1105	83		3		25	60-109
3,3'-Dichlorobenzidine	2667	U	2238	84		2343	88		5		25	64-113
Bis (2-ethylhexyl) phthalate	1333	U	1179	88		1101	83		7		25	50-121
di-n-octyl-phthalate	1333	U	1261	95		1133	85		11		25	56-115
Benzo[b] fluoranthene	1333	U	1148	86		1126	84		2		25	57-114
Benzo[k] fluoranthene	1333	U	1220	91		1131	85		8		25	58-111
Benzo[a] pyrene	1333	U	1068	80		1047	79		2		25	62-111
Indeno [1,2,3-cd] pyrene	1333	U	964	72		967	72		0		25	62-123
Dibenz [a,h] anthrace	1333	U	1017	76		996	75		2		25	69-122
Benzo(g,h,i) perylene	1333	U	1005	75		983	74		2		25	61-122

Column to be used to flag recovery and RPD values with an asterisk

* Values outside of QC limi
RPD: 0 out of 71outside limits
Spike Recovery: 0 out of 142 outside limits

Comments: _____

analytics environmental laboratory LLC

SEMI-VOLATILE ORGANIC SOIL
MATRIX SPIKE/ DUPLICATE
PERCENT RECOVERY

Project #: 218736.002 SDG# 56637

DFTPP File ID: D47794T Site ID: GP Old Town Mill ME ME

Instrument ID: D DFTPP Injection Date:5/31/06

GC Column: RTX-5MS ID: 0.25(mm) DFTPP Injection Time: 1018

Laboratory Control Spike/Duplicate Sample ID: 56637-23

COMPOUND	SPIKE ADDED (ug/kg)	SAMPLE AMT (ug/kg)	MS CONC. (ug/kg)	MS REC %	#	SPIKE ADDED (ug/kg)	MSD CONC. (ug/kg)	MSD REC %	#	RPD	#	QC RPD	LIMITS
Pyridine	1368	U	762	56		1369	840	61		10		25	13-77
N-nitrosodimethylamine	1368	U	892	65		1369	959	70		7		25	20-98
Aniline	1368	U	980	72		1369	977	71		0		25	60-120
Phenol	1368	U	1037	76		1369	1122	82		8		35	51-96
2-Chlorophenol	1368	U	995	73		1369	1052	77		6		50	44-92
bis(2-chloroethyl) ether	1368	U	950	69		1369	1040	76		9		25	36-85
1,3-Dichlorobenzer	1368	U	926	68		1369	984	72		6		25	27-90
1,4-Dichlorobenzer	1368	U	959	70		1369	1020	75		6		27	27-90
1,2-Dichlorobenzer	1368	U	946	69		1369	1038	76		9		25	30-89
Benzyl Alcohol	1368	U	1184	87		1369	1107	81		7		25	45-135
2-Methylphenol	1368	U	1050	77		1369	1090	80		4		25	45-135
bis(2-chloroisopropyl)ether	1368	U	918	67		1369	1003	73		9		25	41-102
3+4-Methylphenol	1368	U	1049	77		1369	1159	85		10		25	45-135
Hexachloroethane	1368	U	842	62		1369	943	69		11		25	30-93
N-nitroso-di-n-propylamine	1368	U	1058	77		1369	1142	83		8		38	46-100
Nitrobenzene	1368	U	1016	74		1369	1099	80		8		25	41-93
Isophorone	1368	U	1166	85		1369	1281	94		9		25	62-116
2-Nitrophenol	1368	U	896	66		1369	992	73		10		25	43-95
2,4-Dimethylphenol	1368	U	949	69		1369	1050	77		10		25	45-79
Bis(2-chloroethoxy)methane	1368	U	969	71		1369	1088	79		12		25	44-100
2,4-Dichlorophenol	1368	U	1102	81		1369	1194	87		8		25	47-95
Benzoic Acid	1368	U	799	58		1369	866	63		8		25	30-150
1,2,4-Trichlorobenzene	1368	U	1070	78		1369	1140	83		6		23	37-92
Naphthalene	1368	U	1037	76		1369	1116	82		7		25	42-99
4-Chloroaniline	1368	U	1110	81		1369	1109	81		0		25	60-120
Hexachlorobutadiene	1368	U	1054	77		1369	1116	82		6		25	35-94
4-Chloro-3-methylphenol	1368	U	1119	82		1369	1232	90		10		33	53-100
2-Methylnaphthalene	1368	U	1176	86		1369	1142	83		3		25	60-120
Hexachlorocyclopentadiene	1368	U	400	29	*	1369	431	32	*	7		25	36-97
2,4,6-Trichlorophenol	1368	U	1160	85		1369	1270	93		9		50	56-100
2,4,5-Trichlorophenol	1368	U	1124	82		1369	1311	96		15		47	45-135
2-Chloronaphthalene	1368	U	1070	78		1369	1168	85		9		25	49-96
2-Nitroaniline	1368	U	1231	90		1369	1276	93		4		25	60-120
2,6-Dinitrotoluene	1368	U	1089	80		1369	1179	86		8		25	60-101
Dimethyl Phthalate	1368	U	1104	81		1369	1215	89		10		25	53-105
Acenaphthylene	1368	U	1177	86		1369	1278	93		8		25	54-97
3-Nitroaniline	1368	U	1321	97		1369	1300	95		2		25	60-120
Acenaphthene	1368	U	1089	80		1369	1202	88		10		19	55-96
2,4-dinitrophenol	1368	U	369	27		1369	372	27		1		25	26-85
Dibenzofuran	1368	U	1234	90		1369	1202	88		3		25	60-120
4-Nitrophenol	1368	U	1401	102		1369	1656	121	*	17		50	57-113

analytics environmental laboratory LLC

SEMI-VOLATILE ORGANIC SOIL
MATRIX SPIKE/ DUPLICATE
PERCENT RECOVERY

Project #: 218736.002 SDG# 56637

DFTPP File ID: D47794T Site ID: GP Old Town Mill ME ME

Instrument ID: D DFTPP Injection Date:5/31/06

GC Column: RTX-5MS ID: 0.25(mm) DFTPP Injection Time: 1018

Laboratory Control Spike/Duplicate Sample ID: 56637-23

COMPOUND	SPIKE ADDED (ug/kg)	SAMPLE AMT (ug/kg)	MS CONC. (ug/kg)	MS REC %	#	SPIKE ADDED (ug/kg)	MSD CONC. (ug/kg)	MSD REC %	#	RPD	#	QC LIMITS RPD	
2,4-Dinitrotoluene	1368	U	1097	80		1369	1181	86		7		47	66-107
Fluorene	1368	U	1119	82		1369	1240	91		10		25	60-101
Diethyl Phthalate	1368	U	1165	85		1369	1255	92		7		25	57-111
4-Chlorophenyl phenyl ether	1368	U	1092	80		1369	1221	89		11		25	54-103
4-Nitroaniline	1368	U	1372	100		1369	1350	99		2		25	60-120
4,6-Dinitro-2-methylphenol	1368	U	270	20	*	1369	294	22	*	9		25	47-110
n-nitrosodiphenylamine	1368	U	1201	88		1369	1322	97		10		25	60-121
Azobenzene	1368	U	997	73		1369	1089	80		9		25	63-115
4-bromophenyl phenyl ether	1368	U	1003	73		1369	1127	82		12		25	55-109
Hexachlorobenzene	1368	U	1090	80		1369	1187	87		9		25	60-104
Pentachlorophenol	1368	U	1188	87		1369	1302	95		9		47	32-115
Phenanthrene	1368	377	1251	64		1369	1362	72		9		25	59-103
Anthracene	1368	U	1190	87		1369	1307	96		9		25	60-107
Carbazole	1368	U	2330	170	*	1369	2571	188	*	10		25	67-113
Di-n-butyl phthalate	1368	U	1083	79		1369	1202	88		10		25	67-119
Fluoranthene	1368	572	1511	69		1369	1647	79		9		25	65-114
Benzidine	2737	U	1196	44		2738	1428	52		18		25	29-187
Pyrene	1368	738	1544	59		1369	1681	69		9		36	52-110
Butyl benzyl phthalate	1368	U	1117	82		1369	1211	89		8		25	56-121
Benzo[a]anthracene	1368	278	1302	75		1369	1444	85		10		25	59-110
Chrysene	1368	349	1376	75		1369	1496	84		8		25	60-109
3,3'-Dichlorobenzidine	2737	U	2263	83		2738	2319	85		2		25	64-113
Bis (2-ethylhexyl) phthalate	1368	U	1091	80		1369	1218	89		11		25	50-121
di-n-octyl-phthalate	1368	U	1951	143	*	1369	2178	159	*	11		25	56-115
Benzo[b] fluoranthene	1368	286	1749	107		1369	1990	125	*	13		25	57-114
Benzo[k] fluoranthene	1368	289	1721	105		1369	1888	117	*	9		25	58-111
Benzo[a] pyrene	1368	364	1308	69		1369	1469	81		12		25	62-111
Indeno [1,2,3-cd] pyrer	1368	U	508	37	*	1369	536	39	*	5		25	62-123
Dibenz [a,h] anthracene	1368	U	573	42	*	1369	611	45	*	7		25	69-122
Benzo(g,h,i) perylene	1368	U	430	31	*	1369	505	37	*	16		25	61-122
	1368												

Column to be used to flag recovery and RPD values with an asterisk

* Values outside of QC limits
RPD: 0 out of 71outside limits
Spike Recovery: 15 out of 142 outside limits

Comments: _____

052406-1

Instrument D

IS only#: S6845 Exp: 11/23/06

Tune File: ABN.u Date: 5/24/06 Loading Check: ARC

Voltage: 1835 Analyst: ARC Unloading Check: SL

Vial	Meth File	Sample Info	Dil	Data File	Win ID	Comments
1	tunemix	S6618 tunemix 50ng	1	D47718T	W-D-052406-1-dup ↑51 Adjust in Tune	
1	↓	S6618 tunemix 50ng	1	19T	New v.cap ↑51 ↓44	
1	↓	S6618 tunemix 50ng	1	20T	Voltage = 1859	
2	S052406	S6852 Pt1 2/4	1	21SI	↓ #47	
Vol	tage = 1894					
1	tunemix	S6618 tunemix 50ng	1	22T	✓	
2	S052406	S6852 Pt1 2/4		23SI	✓	
3		S6851 Pt7A 4/8		24	✓	
4		S6850 Pt2 10/20		25	✓	
5		S6849 Pt3 20/40		26	✓	
6		S6848 Pt4A 40/80		27	✓	
7		S6847 Pt4 80/100		28	✓	
8		S6846 Pt5 100/200		29 ↓	✓	
9		S6857 ABNCC 20/40		30	✓	
10		S6858 Benzidines + Anilines CC 20		31	✓	
1		S6618 tunemix 50ng		32	✓	
5		S6849 Pt3 20/40		33	✓	
11		B052306AS		34	✓	
12		L052306AS		35	✓	
13		U052306AS		36	✓ ↓ being.	
14		~~I06~~ I052306AS-1		37	✓	
15		ARC Spike -2		38	✓	
16		-3		39	✓	
17		-4		40	✓	
18		I052306AW-1		41	✓	
19		-2		42	✓	
20		-3		43	✓	
21		-4		44	✓	
22	↓	56572-2		45	+BS6, -BS6 LL 1:10	

Cont. RS: 22

053006-D

Instrument D

IS only#: S6845 Exp: 11/23/06

Tune File: ABN.u Date: 5/30/06 Loading Check: SC | ALR

Voltage: 1644 Analyst: SC | ALR Unloading Check: SC

Vial	Meth File	Sample Info	Dil	Data File	Win ID	Comments
1	tunemix	S6618 tunemix 50ng	✓	D477SST	✓	W-D-053006-1-ALR
2	S05246	S6849 PT3 20/40	✓	56SC	✓	
3		S6760 Benz + Anil. 20	✓	57SC	✓	
4		B05306AW	✓	58B	✓	
5		L05306AW	✓	59Q	✓	
6		LD05306AW	✓	60Q	a few low	Recoveries (NR)
7		56637-1 8270W	✓	61	↓SS	
8		-2	✓	62	↓SS	
9		-3	✓	63	✓	
10		-20	✓	64	↓SS	
11		-21	✓	65	↓SS	
7		-1, Re-inject	✓	66	Confirms	↓Fs
8		-2, Re-inject	✓	67		
10		-20, Reinjid	✓	68		
11		-21, Re-inject	✓	69		
6	↓	LD05306AW, Re-Inject	✓	70Q	failures	confirm
1	tunemix	S6618 tunemix 50ng	✓	7IT	W-D-053006-2-ALR	
2	S05246	S6845 PT3 20/40	✓	72SC	✓	
3		S6760 Benzidines + Anilines 20	✓	73SC	✓	
4		B05306AS	✓	74B	✓	
5		L05306AS	✓	75Q	✓	
6		LD05306AS	✓	76Q	✓	
7		56633-1 PANS	✓	77	✓	
8		-2 ↓	✓	78	✓	
9		56637-4 8270S	✓	79	✓	
10		-9	✓	80	✓	
11		-10	✓	81	✓	
12		-12	✓	82	✓	
13	↓	-18 ↓	✓	83	✓	

Cont. pg. 25

Cont. 053006-D

Instrument D

IS only#5/6/4 Exp: 11/23/06

Tune File: ABN.u Date: 5/30/06 Loading Check: AMR

Voltage: 1894 Analyst: AMR Unloading Check: SL

Vial	Meth File	Sample Info	Dil	Data File	Win ID	Comments
14	505246	576637-11 82705	/	D47784	✓	
15		-14	/	85	RL 1:2, #IS #6	
16		-19	/	86	✓	
17		-22	/	87	RL 1:2,	#IS #6
18		-23	/	88	RL 1:2,	#IS #6
19		-15	2	RL 1:50 89	500ml +	500ml MeOH
20		-16	5	RL 1:50 90	200ml +	800ml
21	↓	-17	5	RL 1:200 91	↓ +	800ml

5/31/06 SL

L25J

053106-P

Instrument D

IS only#: S6845 Exp: 11/23/06

Tune File: MSN.U Date: 5/31/04 Loading Check: SL | ALR

Voltage: 1844 Analyst: ALR | SL Unloading Check: SL

Vial	Meth File	Sample Info	Dil	Data File	Win ID	Comments
1	tunemix	S6618 tunemix	/	D4772T	W-S-053106-1-ALR ✓	
2	S05246	S6849 PT3 20/40	/	93SL	Need to clip, Failing Compounds	
1	tunemix	S6618 tunemix	/	94T	✓	
2	S05246	S6849 PT3 20/40	/	95SL	✓	
3		S6740 ISTA 20	/	96SL	✓	
4		B05306AS	/	97B	✓	
5		S6637-15, 1:100 82705	100	98	✓	
6		-16, 1:50	50	99	✓	
7		-17, 1:100	100	800	RR 1:20	
8		-14, 1:2	2	801	✓	
9		-22, 1:2	2	802	✓	
10		-23, 1:2	2	803	✓	
11		-17, 1:20	20	804	✓	
12		-23, MS, 1:2	2	805SP	} failures	
13		-23, MSD, 1:2 ↓	2	806SP	} due to Matrix	
14		S6623-1 82705	/	807	✓	
15		S6624-1, 1:2 82705	2	808	RR str	
16		S6572-2, 1:2 PANB	2	809	IS6-31%	
17		S6624-1 82705	/	10	IS6-39%	
1	tunemix	S6862 tunemix 50ng	/	11T	W-S-053106-2-ALR	
2	SIM05316	S6866 SIM Pt1 0.2	/	12SI	✓	
3		S6865 SIM Pt2 1.0	/	13	✓	
4		S6864 SIM Pt3 2.0	/	14	✓	
5		S6863 SIM Pt4 5.0	/	15	✓	
6		S6762 SIM Pt5 10.0	/	16L	✓	
7		S6822 SIM cc 1.0	/	17	✓	
8		B053165W	/	18B	✓	
9		L053165W	/	19Q	✓	
10	↓	LD053165W	/	20Q	✓	

SEMI-VOLATILE ORGANIC
INSTRUMENT PERFORMANCE CHECK
DECAFLUOROTRIPHENYLPHOSPHINE (DFTPP)

Project #:	218736.002	Quote #:	NA	SDG#:	56637
DFTPP File ID:	D47722T.D	Site ID:	GP Old Town Mill	Location:	ME
Instrument ID:	D			DFTPP Injection Date:	5/25/06
GC Column:	RTX-5MS	ID: 0.25(mm)		DFTPP Injection Time:	0305

m/e	ION ABUNDANCE CRITERIA	%RELATIVE ABUNDANCE
51	30 - 60% of mass 198	59.7
68	Less than 2% of mass 69	0.0 (0.0)1
69	Less than 100% of mass 198	70.4
70	Less than 2% of mass 69	0.2 (0.2)1
127	40 - 60% of mass 198	54.1
197	Less than 1% of mass 198	0.0
198	Base peak, 100% relative abundance	100.0
199	5 - 9% of mass 198	6.6
275	10 - 30% of mass 198	25.5
365	1 - 100% of mass 198	3.3
441	.01 - 100% of mass 443	6.4 (42.0)2
442	40 - 100% of mass 198	83.1
443	17 - 23% of mass 442	15.3 (18.4)3

1-Value is % mass 692-Value is % mass 4 3-Value is % mass 442

This check applies to the following SAMPLES, MS, MSD, BLANKS and STANDARDS:

	SAMPLE NO.	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED	TIME ANALYZED
01	S6852 PT1 2/4	S6852 PT1 2/4	D47723SI.D	5/25/06	3:47
02	S6851 PT2A 4/8	S6851 PT2A 4/8	D47724SI.D	5/25/06	4:20
03	S6850 PT2 10/20	S6850 PT2 10/20	D47725SI.D	5/25/06	4:53
04	S6849 PT3 20/40	S6849 PT3 20/40	D47726SI.D	5/25/06	5:26
05	S6848 PT4A 40/80	S6848 PT4A 40/80	D47727SI.D	5/25/06	5:59
06	S6847 PT4 50/100	S6847 PT4 50/100	D47728SI.D	5/25/06	6:32
07	S6846 PT5 100/200	S6846 PT5 100/200	D47729SI.D	5/25/06	7:06
08	S6857 ABN CC 20/40	S6857 ABN CC 20	D47730.[	5/25/06	7:39
09	S6858 BENZIDINES +	S6858 BENZIDINE	D47731.[	5/25/06	8:12
10					
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SEMI-VOLATILE ORGANIC
INSTRUMENT PERFORMANCE CHECK
DECAFLUOROTRIPHENYLPHOSPHINE (DFTPP)

Project #: 218736.002 Quote #: NA SDG#: 56637

DFTPP File ID: D47755T.D Site ID: GP Old Town Mill Location: ME

Instrument ID: D DFTPP Injection Date: 5/30/06

GC Column: RTX-5MS ID: 0.25(mm) DFTPP Injection Time: 0639

m/e	ION ABUNDANCE CRITERIA	%RELATIVE ABUNDANCE
51	30 - 60% of mass 198	44.9
68	Less than 2% of mass 69	0.0 (0.0)1
69	Less than 100% of mass 198	56.5
70	Less than 2% of mass 69	0.3 (0.5)1
127	40 - 60% of mass 198	47.5
197	Less than 1% of mass 198	0.0
198	Base peak, 100% relative abunda	100.0
199	5 - 9% of mass 198	6.7
275	10 - 30% of mass 198	26.5
365	1 - 100% of mass 198	3.6
441	.01 - 100% of mass 443	13.9 (82.7)2
442	40 - 100% of mass 198	94.1
443	17 - 23% of mass 442	16.8 (17.9)3

1-Value is % mass 69 2-Value is % mass 43 3-Value is % mass 442

This check applies to the following SAMPLES, MS, MSD, BLANKS and STANDARDS:

	SAMPLE NO.	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED	TIME ANALYZED
01	S6849 PT3 20/40	S6849 PT3 20/40	D47756SC.	5/30/06	7:24
02	S6760 B+A 20	S6760 B+A 20	D47757SC.	5/30/06	7:57
03	B05306AW	B05306AW	D47758B.D	5/30/06	9:36
04	L05306AW	L05306AW	D47759Q.D	5/30/06	10:09
05	LD05306AW	LD05306AW	D47760Q.D	5/30/06	10:42
06	56637-1	56637-1	D47761.I	5/30/06	10:56
07	56637-2	56637-2	D47762.I	5/30/06	12:02
08	56637-3	56637-3	D47763.I	5/30/06	12:35
09	56637-20	56637-2(	D47764.I	5/30/06	13:08
10	56637-21	56637-2	D47765.I	5/30/06	13:50
11	56637-1, RE-INJECT	56637-1, RE-INJE	D47766.I	5/30/06	14:11
12	56637-2, RE-INJECT	56637-2, RE-INJE	D47767.I	5/30/06	14:44
13	56637-20, RE-INJECT	56637-20, RE-INJ	D47768.I	5/30/06	15:17
14	56637-21, RE-INJECT	56637-21, RE-INJ	D47769.I	5/30/06	15:50
15	LD05306AW, RE-INJEC	LD05306AW, RE-I	D47770Q.D	5/30/06	16:23
16					
17					
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SEMI-VOLATILE ORGANIC
INSTRUMENT PERFORMANCE CHECK
DECAFLUOROTRIPHENYLPHOSPHINE (DFTPP)

Project #: 218736.002 Quote #: NA SDG#: 56637

DFTPP File ID: D47771T.D Site ID: GP Old Town Mill Location: ME

Instrument ID: D DFTPP Injection Date: 5/30/06

GC Column: RTX-5MS ID: 0.25(mm) DFTPP Injection Time: 2043

m/e	ION ABUNDANCE CRITERIA	%RELATIVE ABUNDANCE
51	30 - 60% of mass 198	57.6
68	Less than 2% of mass 69	0.0 (0.0)1
69	Less than 100% of mass 198	67.3
70	Less than 2% of mass 69	0.2 (0.2)1
127	40 - 60% of mass 198	54.0
197	Less than 1% of mass 198	0.0
198	Base peak, 100% relative abundance	100.0
199	5 - 9% of mass 198	7.1
275	10 - 30% of mass 198	23.4
365	1 - 100% of mass 198	2.6
441	.01 - 100% of mass 443	9.4 (78.0)2
442	40 - 100% of mass 198	62.1
443	17 - 23% of mass 442	12.1 (19.5)3

1-Value is % mass 692-Value is % mass 4 3-Value is % mass 442

This check applies to the following SAMPLES, MS, MSD, BLANKS and STANDARDS:

	SAMPLE NC	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED	TIME ANALYZED
01	S6849 PT3 20/40	S6849 PT3 20/40	D47772SC.	5/30/06	21:08
02	S6760 B+A 2	S6760 B+A 20	D47773SC.	5/30/06	21:46
03	B05306AS	B05306AS	D47774B.D	5/30/06	22:19
04	L05306AS	L05306AS	D47775Q.D	5/30/06	23:02
05	LD05306AS	LD05306AS	D47776Q.D	5/30/06	23:17
06	ZZZZZ	ZZZZZ	D47777.I	5/30/06	23:50
07	ZZZZZ	ZZZZZ	D47778.I	5/31/06	0:23
08	56637-4	56637-4	D47779.I	5/31/06	1:23
09	56637-9	56637-9	D47780.I	5/31/06	1:58
10	56637-1	56637-1	D47781.I	5/31/06	2:31
11	56637-1	56637-1	D47782.I	5/31/06	3:04
12	56637-1	56637-1	D47783.I	5/31/06	3:37
13	56637-11	56637-11	D47784.I	5/31/06	4:10
14	56637-1	56637-1	D47785.I	5/31/06	4:43
15	56637-1	56637-1	D47786.I	5/31/06	5:16
16	56637-2	56637-2	D47787.I	5/31/06	5:49
17	56637-2	56637-2	D47788.I	5/31/06	6:22
18	56637-15,,1:2	56637-15,,1:2	D47789.I	5/31/06	6:55
19	56637-16,,1:5	56637-16,,1:5	D47790.I	5/31/06	7:28
20	56637-17,,1:5	56637-17,,1:5	D47791.I	5/31/06	8:01
21					
22					
23					

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SEMI-VOLATILE ORGANIC
INSTRUMENT PERFORMANCE CHECK
DECAFLUOROTRIPHENYLPHOSPHINE (DFTPP)

Project #: 218736.002 Quote #: NA SDG#: 56637

DFTPP File ID: D47794T. Site ID: GP Old Town Mill Location: ME

Instrument ID: D DFTPP Injection Date: 5/31/06

GC Column: RTX-5MS ID: 0.25(mm) DFTPP Injection Time: 1018

m/e	ION ABUNDANCE CRITERIA	%RELATIVE ABUNDANCE	
51	30 - 60% of mass 198	45.6	
68	Less than 2% of mass 69	0.0 (	0.0)1
69	Less than 100% of mass 198	53.1	
70	Less than 2% of mass 69	0.1 (	0.2)1
127	40 - 60% of mass 198	48.6	
197	Less than 1% of mass 198	0.0	
198	Base peak, 100% relative ab	100.0	
199	5 - 9% of mass 198	6.3	
275	10 - 30% of mass 198	25.7	
365	1 - 100% of mass 198	3.4	
441	.01 - 100% of mass 443	13.7 (	85.2)2
442	40 - 100% of mass 198	89.1	
443	17 - 23% of mass 442	16.1 (	18.1)3

1-Value is % mass 69 2-Value is % mass 4433-Value is % mass 442

This check applies to the following SAMPLES, MS, MSD, BLANKS and STANDARDS:

	SAMPLE NO.	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED	TIME ANALYZED
01	S6849 PT3 20/40	S6849 PT3 20/40	D47795SC.D	5/31/06	11:03
02	S6740 B+A 20	S6740 B+A	D47796SC.D	5/31/06	11:40
03	B05306AS	B05306AS	D47797B.D	5/31/06	12:13
04	56637-15,,1:100	56637-15,,1:100	D47798.D	5/31/06	12:46
05	56637-16,,1:50	56637-16,,1:50	D47799.D	5/31/06	13:19
06	56637-17,,1:100	56637-17,,1:100	D47800.D	5/31/06	14:07
07	56637-14,DL,1:2	56637-14,DL,1:2	D47801.D	5/31/06	14:59
08	56637-22,DL,1:2	56637-22,DL,1:2	D47802.D	5/31/06	15:32
09	56637-23,DL,1:2	56637-23,DL,1:2	D47803.D	5/31/06	16:10
10	56637-17,,1:20	56637-17,,1:20	D47804.D	5/31/06	16:43
11	56637-23.MS,,1:2	56637-23.MS,,1:2	D47805SP.D	5/31/06	17:16
12	56637-23.MSD,,1:2	56637-23.MSD,,1:2	D47806SP.D	5/31/06	17:49
13	ZZZZZ	ZZZZZ	D47807.D	5/31/06	18:22
14	ZZZZZ	ZZZZZ	D47808.D	5/31/06	18:55
15	ZZZZZ	ZZZZZ	D47809.D	5/31/06	19:28
16	ZZZZZ	ZZZZZ	D47810.D	5/31/06	20:35
17					
18					
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20					
21					
22					
23					

Response Factor Report D Instrum

Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Fri May 26 09:35:43 2006
Response via : Initial Calibration

5.26.06

5.26.06 SL

Calibration Files
1 =D47723SI.D 2 =D47725SI.D 3 =D47726SI.D
4 =D47728SI.D 5 =D47729SI.D 2A =D47724SI.D

		Compound	1	2	3	4	5	2A	Avg	%RSD
1)	I	d4-1,4-dichlorobenzen	ISTD							
2)	TM	Pyridine	1.297	1.808	1.592	1.625	1.589	1.440	1.570	10.29
3)	TM	N-nitrosodimethylam	1.410	1.169	1.267	1.203	1.176	1.167	1.230	7.07
4)	S	2-Fluorophenol	1.454	1.520	1.567	1.577	1.546	1.520	1.538	2.89
5)	S	d5-Phenol	1.735	1.801	1.872	1.903	1.929	1.770	1.844	4.08
6)	TM	Aniline	2.292	2.407	2.506	2.473	2.548	2.389	2.438	3.47
7)	TMC	Phenol	1.982	1.936	2.003	2.054	2.064	1.971	2.004	2.28
8)	TM	2-Chlorophenol	1.523	1.511	1.540	1.575	1.577	1.481	1.537	2.26
9)	TM	bis(2-chloroethyl)	1.629	1.634	1.623	1.721	1.516	1.490	1.616	5.32
10)	TM	1,3-Dichlorobenzene	1.486	1.449	1.493	1.517	1.473	1.506	1.489	1.53
11)	TMC	1,4-Dichlorobenzene	1.487	1.513	1.580	1.553	1.531	1.499	1.532	2.26
12)	TM	1,2-Dichlorobenzene	1.431	1.420	1.472	1.464	1.436	1.413	1.441	1.54
13)	TM	Benzyl Alcohol	0.917	1.022	1.115	1.112	1.116	1.000	1.056	7.42
14)	TM	2-Methylphenol	1.366	1.424	1.456	1.460	1.447	1.396	1.428	2.48
15)	TM	bis(2-chloroisoprop	2.381	2.481	2.514	2.497	2.348	2.482	2.451	2.54
16)	TM	3+4-Methylphenol	1.364	1.448	1.528	1.542	1.547	1.444	1.485	4.59
17)	TM	Hexachloroethane	0.715	0.710	0.723	0.735	0.715	0.706	0.718	1.31
18)	TMP	N-nitroso-di-n-prop	0.975	0.948	0.962	0.974	0.956	0.972	0.966	1.15
19)	I	d8-Naphthalene	ISTD							
20)	S	d5-nitrobenzene	0.391	0.450	0.458	0.454	0.443	0.421	0.439	5.58
21)	TM	Nitrobenzene	0.202	0.225	0.233	0.231	0.232	0.218	0.225	5.06
22)	TM	Isophorone	0.712	0.737	0.753	0.738	0.720	0.720	0.732	2.06
23)	TCM	2-Nitrophenol	0.193	0.224	0.228	0.238	0.236	0.215	0.225	7.16
24)	TM	2,4-Dimethylphenol	0.368	0.418	0.422	0.423	0.420	0.406	0.412	4.88
25)	TM	Bis(2-chloroethoxy)	0.451	0.501	0.501	0.506	0.487	0.473	0.489	4.18
26)	TCM	2,4-Dichlorophenol	0.234	0.285	0.290	0.298	0.301	0.277	0.283	8.20
27)	TM	Benzoic Acid	0.072	0.187	0.235	0.292	0.309	0.129	0.215	~~41.79~~ Q
28)	TM	1,2,4-Trichlorobenz	0.282	0.301	0.303	0.309	0.309	0.303	0.302	3.12
29)	TM	Naphthalene	1.052	1.098	1.109	1.102	1.073	1.077	1.088	2.01
30)	T	2,6-Dichlorophenol	0.273	0.286	0.291	0.298	0.317	0.277	0.291	5.05
31)	TM	4-Chloroaniline	0.453	0.478	0.489	0.508	0.540	0.452	0.488	6.39
32)	TCM	Hexachlorobutadiene	0.161	0.164	0.169	0.170	0.172	0.166	0.167	2.21
33)	TMC	4-Chloro-3-methylph	0.299	0.332	0.354	0.366	0.384	0.324	0.345	8.29
34)	TM	2-Methylnaphthalene	0.558	0.635	0.647	0.652	0.698	0.601	0.633	6.91
35)	I	d10-Acenaphthene	ISTD							
36)	TPM	Hexachlorocyclopent	0.053	0.108	0.173	0.220	0.016	0.122	~~63.44~~ Q	
37)	TCM	2,4,6-Trichlorophen	0.318	0.368	0.383	0.386	0.403	0.343	0.371	8.24
38)	TM	2,4,5-Trichlorophen	0.337	0.352	0.385	0.396	0.511	0.354	0.393	14.91
39)	S	2-Fluorobiphenyl	1.304	1.320	1.341	1.326	1.444	1.302	1.342	3.64
40)	TM	2-Chloronaphthalene	1.129	1.123	1.174	1.187	1.362	1.155	1.191	6.83
41)	TM	2-Nitroaniline	0.405	0.502	0.531	0.527	0.530	0.479	0.502	9.56

(#) = Out of Range ### Number of calibration levels exceeded format ###

Response Factor Report D Instrum

Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Fri May 26 09:35:43 2006
Response via : Initial Calibration

Calibration Files
1 =D47723SI.D 2 =D47725SI.D 3 =D47726SI.D
4 =D47728SI.D 5 =D47729SI.D 2A =D47724SI.D

		Compound	1	2	3	4	5	2A	Avg	%RSD	
42)	TM	2,6-Dinitrotoluene	0.315	0.340	0.355	0.362	0.357	0.319	0.345	6.19	
43)	TM	Dimethyl Phthalate	1.356	1.371	1.404	1.399	1.429	1.375	1.396	2.25	
44)	TM	Acenaphthylene	1.788	1.842	1.906	1.872	1.872	1.819	1.860	2.49	
45)	TM	3-Nitroaniline	0.310	0.383	0.412	0.427	0.430	0.389	0.397	10.95	
46)	TMC	Acenaphthene	1.017	1.037	1.055	1.078	1.093	1.008	1.057	3.60	
47)	TPM	2,4-dinitrophenol	0.050	0.139	0.188	0.229	0.253	0.082	0.167	~~46.92~~	Q
48)	TM	Dibenzofuran	1.674	1.655	1.652	1.645	1.632	1.678	1.661	1.24	
49)	TMP	4-Nitrophenol	0.165	0.146	0.155	0.170	0.180	0.172	0.165	6.86	
50)	TM	2,4-Dinitrotoluene	0.347	0.449	0.472	0.476	0.482	0.436	0.450	10.83	
51)	T	2,3,4,6-Tetrachloro	0.170	0.239	0.259	0.278	0.302	0.222	0.250	~~17.89~~	Q
52)	TM	Fluorene	1.262	1.288	1.301	1.302	1.300	1.280	1.296	1.87	
53)	TM	Diethyl Phthalate	1.366	1.375	1.421	1.411	1.221	1.432	1.380	5.46	
54)	TM	4-Chlorophenyl phen	0.596	0.584	0.613	0.607	0.669	0.582	0.610	4.85	
55)	TM	4-Nitroaniline	0.377	0.409	0.432	0.414	0.349	0.404	0.403	7.59	
56)	TM	4,6-Dinitro-2-methy	0.156	0.265	0.297	0.316	0.308	0.201	0.266	~~24.07~~	Q
57)	TCM	n-nitrosodiphenylam	0.835	0.857	0.878	0.867	0.942	0.864	0.876	3.89	
58)	TM	Azobenzene	1.861	1.773	1.938	1.894	1.927	1.908	1.897	3.47	
59)	TM	Biphenyl	1.407	1.465	1.486	1.502	1.732	1.455	1.513	6.96	
60)	S	2,4,6-Tribromopheno	0.187	0.205	0.215	0.224	0.238	0.198	0.214	8.44	
61)	I	d10-Phenanthrene	ISTD								
62)	TM	4-bromophenyl pheny	0.206	0.212	0.214	0.224	0.223	0.212	0.216	3.19	
63)	TM	Hexachlorobenzene	0.220	0.243	0.260	0.258	0.266	0.247	0.251	6.17	
64)	TMC	Pentachlorophenol		0.091	0.111	0.140	0.151	0.066	0.117	28.45#	Q
65)	TM	Phenanthrene	0.975	1.005	1.000	1.004	1.060	0.992	1.009	2.66	
66)	TM	Anthracene	1.034	1.039	1.055	1.021	0.938	1.009	1.018	3.71	
67)	TM	Carbazole	0.981	0.990	0.997	0.971	0.935	0.996	0.984	2.62	
68)	TM	Di-n-butyl phthalat	1.366	1.382	1.441	1.413	1.159	1.373	1.367	7.08	
69)	TCM	Fluoranthene	0.995	1.064	1.063	1.080	0.953	1.014	1.039	5.13	
70)	TM	Benzidine	0.566	0.573	0.541	0.464	0.334	0.587	0.507	17.58	
71)	I	d12-Chrysene	ISTD								
72)	TM	Pyrene	1.196	1.200	1.215	1.218	1.192	1.178	1.204	1.48	
73)	S	d14-p-terphenyl	0.790	0.833	0.849	0.871	0.953	0.824	0.855	5.97	
74)	TM	Butyl benzyl phthal	0.667	0.708	0.733	0.729	0.736	0.705	0.717	3.58	
75)	TM	Benzo[a]anthracene	1.060	1.075	1.080	1.083	1.058	1.059	1.071	1.04	
76)	TM	Chrysene	0.974	1.004	1.026	1.022	1.041	0.984	1.011	2.42	
77)	TM	3,3'-Dichlorobenzid	0.416	0.435	0.458	0.473	0.496	0.433	0.455	6.17	
78)	TM	Bis (2-ethylhexyl)	0.919	0.965	0.994	1.020	1.135	0.950	1.000	6.98	
79)	I	d12-Perylene	ISTD								
80)	TCM	di-n-octyl-phthalat	1.600	1.760	1.805	1.830	1.796	1.703	1.764	4.92	
81)	TM	Benzo[b] fluoranthe	1.098	1.082	1.112	1.126	1.098	1.049	1.103	3.15	
82)	TM	Benzo[k] fluoranthe	1.020	1.074	1.087	1.080	1.041	1.051	1.063	2.44	

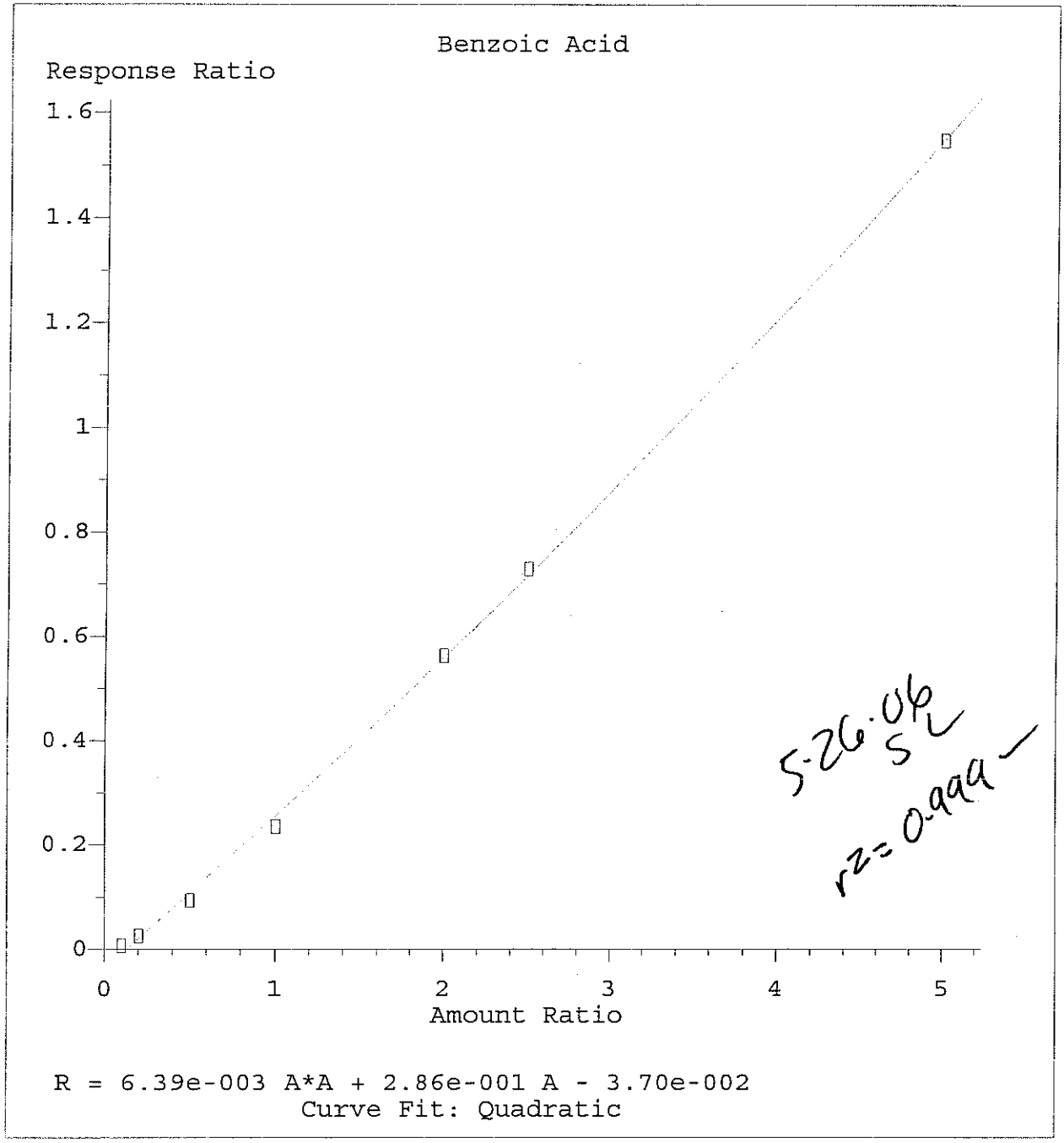
(#) = Out of Range ### Number of calibration levels exceeded format ###

Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Fri May 26 09:35:43 2006
Response via : Initial Calibration

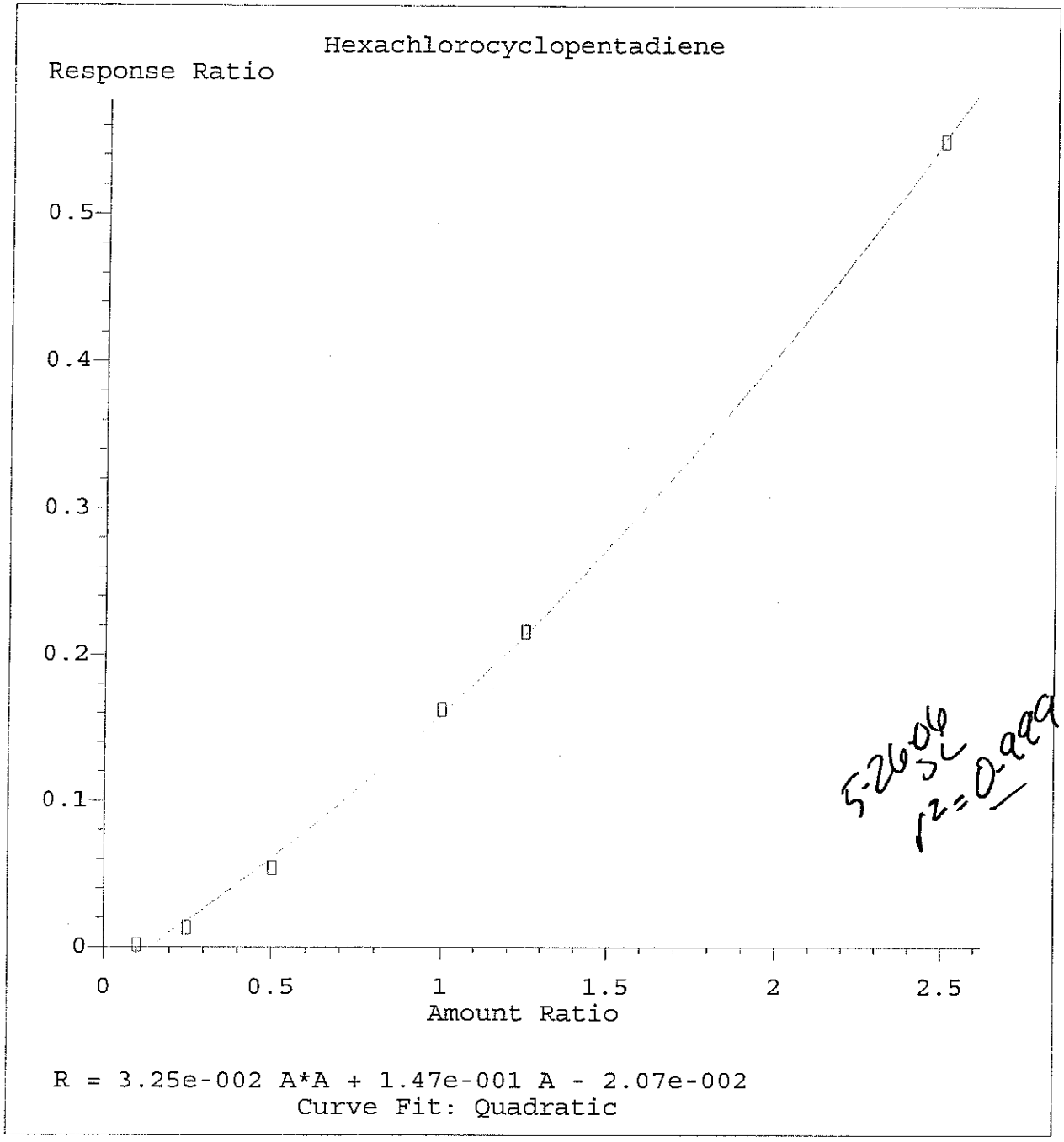
Calibration Files
1 =D47723SI.D 2 =D47725SI.D 3 =D47726SI.D
4 =D47728SI.D 5 =D47729SI.D 2A =D47724SI.D

	Compound	1	2	3	4	5	2A	Avg	%RSD
83)	TCM Benzo[a] pyrene	0.986	1.033	1.060	1.064	1.037	1.034	1.042	2.91
84)	TM Indeno [1,2,3-cd] p	0.854	0.969	1.021	1.118	1.191	0.918	1.023	11.58
85)	TM Dibenz [a,h] anthra	0.821	0.975	1.003	1.076	1.175	0.886	0.998	11.88
86)	TM Benzo(g,h,i) peryl	0.865	0.984	1.010	1.065	1.110	0.938	1.002	8.22

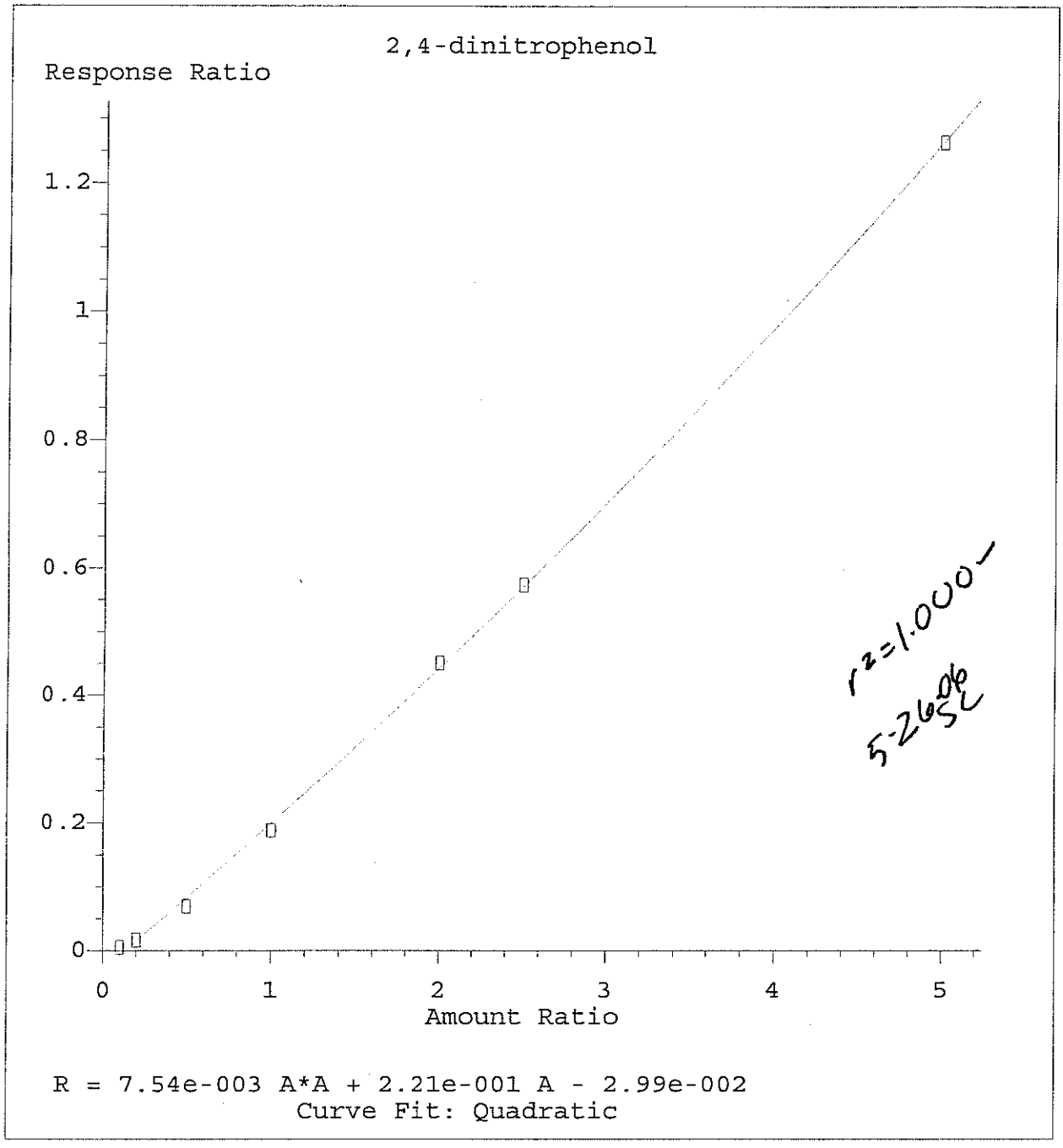
(#) = Out of Range ### Number of calibration levels exceeded format ###



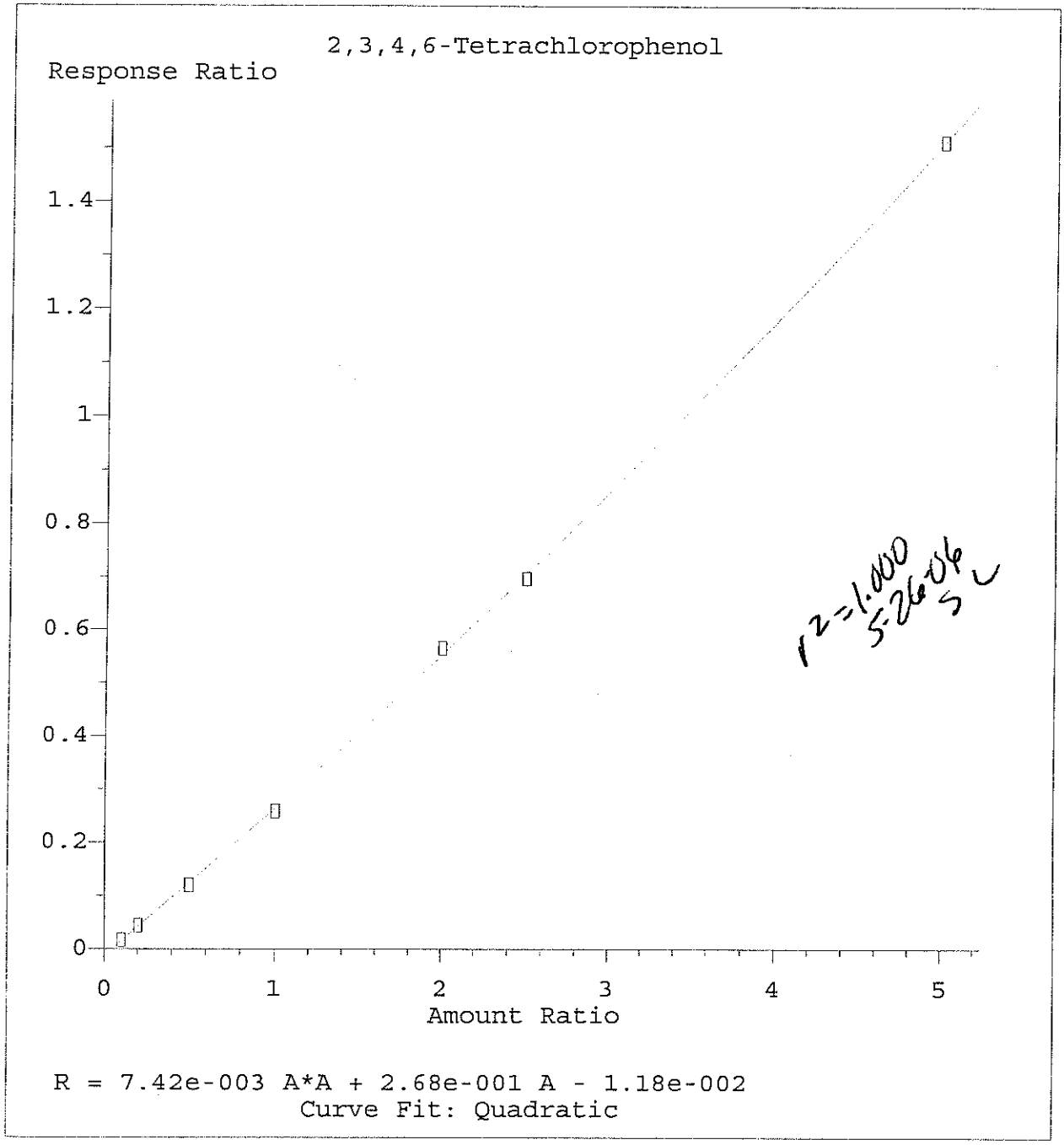
Method Name: D:\HPCHEM\1\METHODS\S05246.M
Calibration Table Last Updated: Fri May 26 09:33:08 2006



Method Name: D:\HPCHEM\1\METHODS\S05246.M
Calibration Table Last Updated: Fri May 26 09:33:30 2006

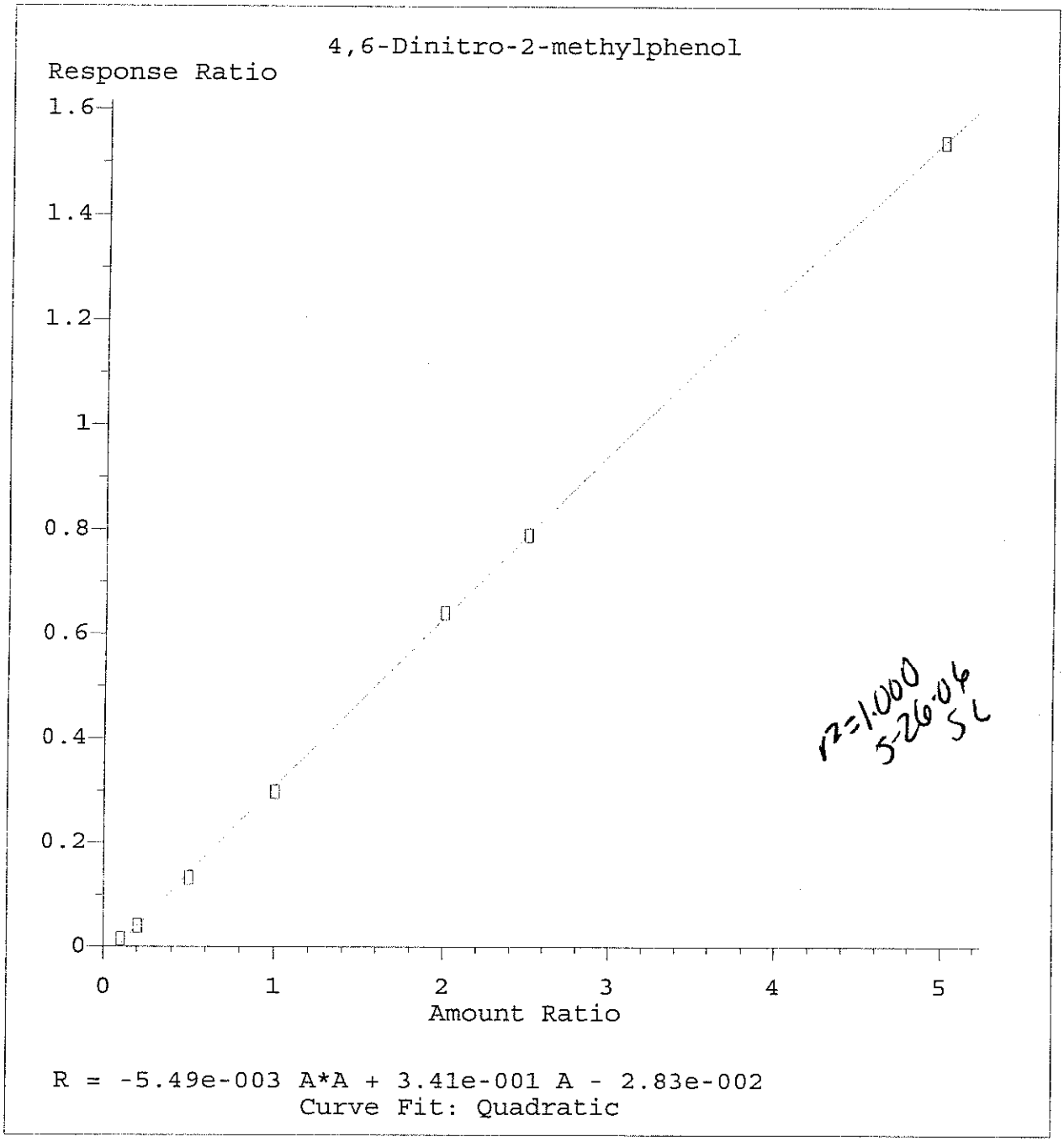


Method Name: D:\HPCHEM\1\METHODS\S05246.M
Calibration Table Last Updated: Fri May 26 09:33:57 2006

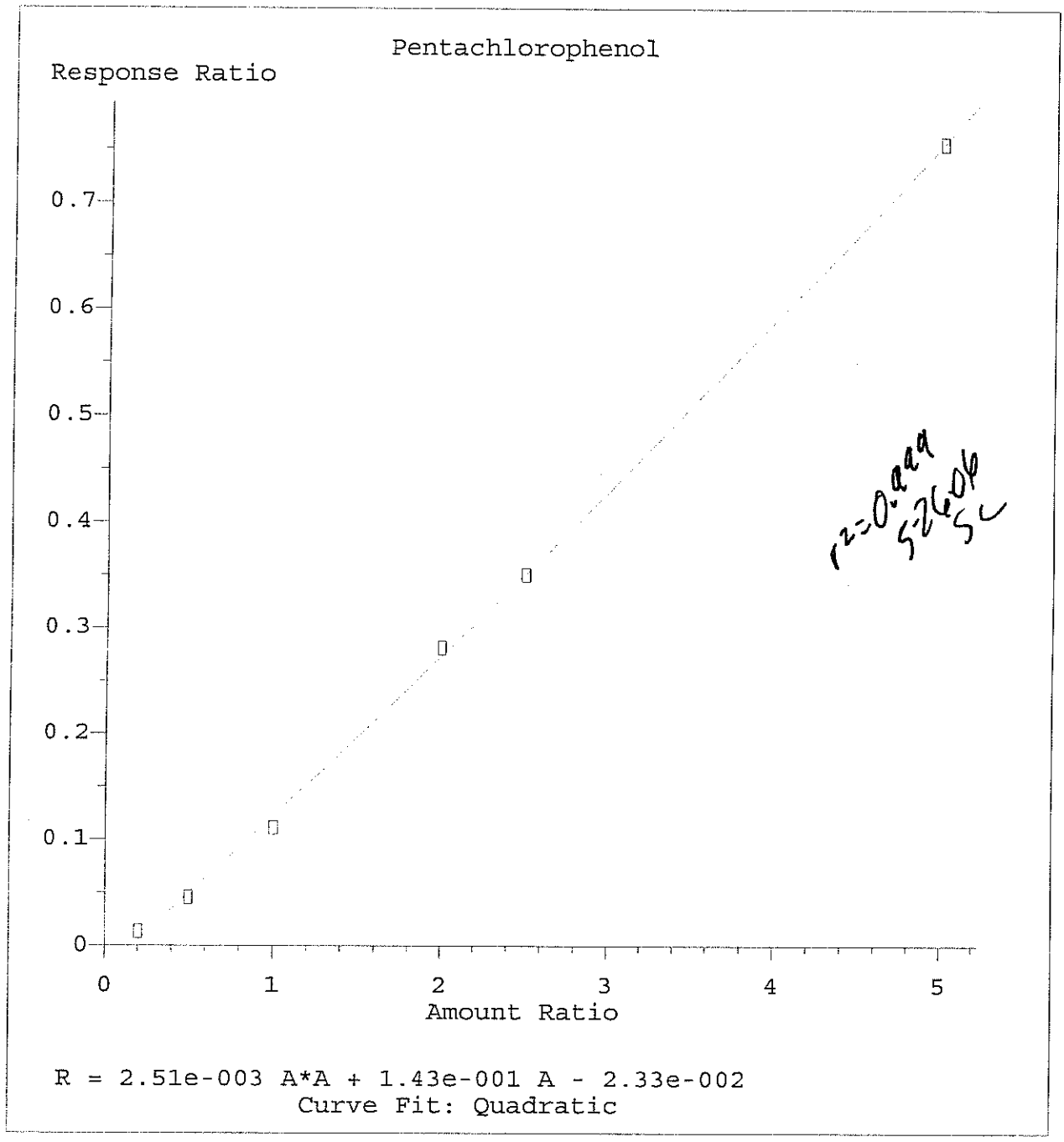


$r^2 = 1.000$
5-26-06
SL

Method Name: D:\HPCHEM\1\METHODS\S05246.M
Calibration Table Last Updated: Fri May 26 09:34:20 2006



Method Name: D:\HPCHEM\1\METHODS\S05246.M
Calibration Table Last Updated: Fri May 26 09:35:07 2006



Method Name: D:\HPCHEM\1\METHODS\S05246.M
Calibration Table Last Updated: Fri May 26 09:35:22 2006

Quantitation Report (QT Reviewed)

Data File : D:\HPCHEM\1\DATA\052406-D\D47723SI.D Vial: 2
Acq On : 25 May 2006 3:47 am Operator:
Sample : S6852 PT1 2/4 Inst : D Instrum
Misc : Multiplr: 1.00
Integration Params: RTEINT.P
Quant Time: May 26 9:23 2006 Quant Results File: S05246.RES

Quant Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Fri May 26 09:21:45 2006
Response via : Initial Calibration
DataAcq Meth : ABN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) d4-1,4-dichlorobenzene	5.04	152	499606	40.00	ug/ml	0.00
19) d8-Naphthalene	7.95	136	1884971	40.00	ug/ml	0.00
35) d10-Acenaphthene	12.21	164	940184	40.00	ug/ml	0.00
61) d10-Phenanthrene	15.75	188	1589687	40.00	ug/ml	0.00
71) d12-Chrysene	22.17	240	1403454	40.00	ug/ml	0.00
79) d12-Perylene	25.80	264	1349553	40.00	ug/ml	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	2.73	112	72662	4.07	ug/ml	0.00
Spiked Amount 100.000	Range 28 - 110		Recovery = 4.07%#			
5) d5-Phenol	4.62	99	86660	4.10	ug/ml	0.02
Spiked Amount 100.000	Range 22 - 110		Recovery = 4.10%#			
20) d5-nitrobenzene	6.36	82	73648	3.45	ug/ml	-0.01
Spiked Amount 50.000	Range 45 - 107		Recovery = 6.90%#			
39) 2-Fluorobiphenyl	10.63	172	122592	3.94	ug/ml	-0.01
Spiked Amount 50.000	Range 44 - 102		Recovery = 7.88%#			
60) 2,4,6-Tribromophenol	14.18	330	17569	2.90	ug/ml	0.00
Spiked Amount 100.000	Range 63 - 127		Recovery = 2.90%#			
73) d14-p-terphenyl	19.69	244	110876	3.67	ug/ml	-0.02
Spiked Amount 50.000	Range 33 - 132		Recovery = 7.34%#			
Target Compounds						Qvalue
2) Pyridine	1.03	79	32390	1.19	ug/ml#	100
3) N-nitrosodimethylamine	1.00	42	35225	2.24	ug/ml	82
6) Aniline	4.47	93	57251	1.95	ug/ml	96
7) Phenol	4.65	94	99011	4.31	ug/ml	95
8) 2-Chlorophenol	4.74	128	76113	4.01	ug/ml	97
9) bis(2-chloroethyl) ether	4.59	93	40682	2.11	ug/ml#	79
10) 1,3-Dichlorobenzene	4.96	146	37121	1.86	ug/ml	96
11) 1,4-Dichlorobenzene	5.07	146	37147	1.84	ug/ml#	69
12) 1,2-Dichlorobenzene	5.47	146	35753	1.88	ug/ml	99
13) Benzyl Alcohol	5.47	108	22912	2.04	ug/ml	89
14) 2-Methylphenol	5.88	108	68229	4.31	ug/ml	97
15) bis(2-chloroisopropyl)ethe	5.77	45	59490	1.59	ug/ml	99
16) 3+4-Methylphenol	6.23	108	68144	4.19	ug/ml	99
17) Hexachloroethane	6.13	117	17867	2.14	ug/ml	98
18) N-nitroso-di-n-propylamine	6.10	70	24352	1.59	ug/ml	97
21) Nitrobenzene	6.40	123	19041	1.88	ug/ml	81
22) Isophorone	6.93	82	67066	1.81	ug/ml	98
23) 2-Nitrophenol	7.15	139	36470	3.25	ug/ml	93
24) 2,4-Dimethylphenol	7.40	107	69442	4.19	ug/ml	96
25) Bis(2-chloroethoxy)methane	7.54	93	42511	1.93	ug/ml	98

(#) = qualifier out of range (m) = manual integration

Quantitation Report (QT Reviewed)

Data File : D:\HPCHEM\1\DATA\052406-D\D47723SI.D Vial: 2
Acq On : 25 May 2006 3:47 am Operator:
Sample : S6852 PT1 2/4 Inst : D Instrum
Misc : Multiplr: 1.00
Integration Params: RTEINT.P
Quant Time: May 26 9:23 2006 Quant Results File: S05246.RES

Quant Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Fri May 26 09:21:45 2006
Response via : Initial Calibration
DataAcq Meth : ABN

Compound	R.T.	QIon	Response	Conc Unit	Qvalue
26) 2,4-Dichlorophenol	7.82	162	44114	3.38 ug/ml	94
27) Benzoic Acid	7.75	105	13546m	3.97 ug/ml	
28) 1,2,4-Trichlorobenzene	7.87	180	26584	1.89 ug/ml	97
29) Naphthalene	7.99	128	99128	1.96 ug/ml	98
30) 2,6-Dichlorophenol	8.28	162	51428	3.93 ug/ml	99
31) 4-Chloroaniline	8.27	127	42721	1.90 ug/ml	96
32) Hexachlorobutadiene	8.45	225	15185	1.85 ug/ml	97
33) 4-Chloro-3-methylphenol	9.67	107	56298	4.11 ug/ml	96
34) 2-Methylnaphthalene	9.66	141	52594	1.83 ug/ml	96
36) Hexachlorocyclopentadiene	10.20	237	61	3.89 ug/ml#	21
37) 2,4,6-Trichlorophenol	10.51	196	29929	3.07 ug/ml	93
38) 2,4,5-Trichlorophenol	10.73	196	31723	3.16 ug/ml	95
40) 2-Chloronaphthalene	10.81	162	53069	1.84 ug/ml	96
41) 2-Nitroaniline	11.26	138	19024	1.38 ug/ml	96
42) 2,6-Dinitrotoluene	11.95	165	14813	1.71 ug/ml	100
43) Dimethyl Phthalate	11.79	163	63729	1.87 ug/ml	97
44) Acenaphthylene	11.82	152	84051	1.75 ug/ml	99
45) 3-Nitroaniline	12.28	138	14556	1.34 ug/ml	93
46) Acenaphthene	12.28	154	47800	1.79 ug/ml	99
47) 2,4-dinitrophenol	12.59	184	4691m	5.80 ug/ml	
48) Dibenzofuran	12.67	168	78685	1.94 ug/ml	95
49) 4-Nitrophenol	13.16	109	15483m	4.04 ug/ml	
50) 2,4-Dinitrotoluene	12.89	165	16305	1.48 ug/ml#	80
51) 2,3,4,6-Tetrachlorophenol	13.18	232	15958	2.11 ug/ml	98
52) Fluorene	13.50	166	59303	1.91 ug/ml	99
53) Diethyl Phthalate	13.50	149	64204	1.78 ug/ml	99
54) 4-Chlorophenyl phenyl eth	13.56	204	28029	1.79 ug/ml#	88
55) 4-Nitroaniline	13.77	138	17724	1.58 ug/ml	93
56) 4,6-Dinitro-2-methylphenol	13.87	198	14621	4.93 ug/ml	99
57) n-nitrosodiphenylamine	13.90	169	39240	1.86 ug/ml	99
58) Azobenzene	13.93	77	87488	1.78 ug/ml#	87
59) Biphenyl	10.83	154	66154	1.76 ug/ml	98
62) 4-bromophenyl phenyl ether	14.72	248	16376	1.70 ug/ml	98
63) Hexachlorobenzene	15.02	284	17524	1.60 ug/ml	90
64) Pentachlorophenol	15.56	266	7480m	1.30 ug/ml	
65) Phenanthrene	15.79	178	77503	1.85 ug/ml	98
66) Anthracene	15.90	178	82188	1.95 ug/ml	99
67) Carbazole	16.40	167	77999	1.92 ug/ml	97
68) Di-n-butyl phthalate	17.44	149	108615	1.79 ug/ml#	97
69) Fluoranthene	18.67	202	79126	1.75 ug/ml	99
70) Benzidine	19.12	184	44989	2.36 ug/ml#	97

(#) = qualifier out of range (m) = manual integration

Quantitation Report (QT Reviewed)

Data File : D:\HPCHEM\1\DATA\052406-D\D47723SI.D Vial: 2
Acq On : 25 May 2006 3:47 am Operator:
Sample : S6852 PT1 2/4 Inst : D Instrum
Misc : Multiplr: 1.00
Integration Params: RTEINT.P
Quant Time: May 26 9:23 2006 Quant Results File: S05246.RES

Quant Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Fri May 26 09:21:45 2006
Response via : Initial Calibration
DataAcq Meth : ABN

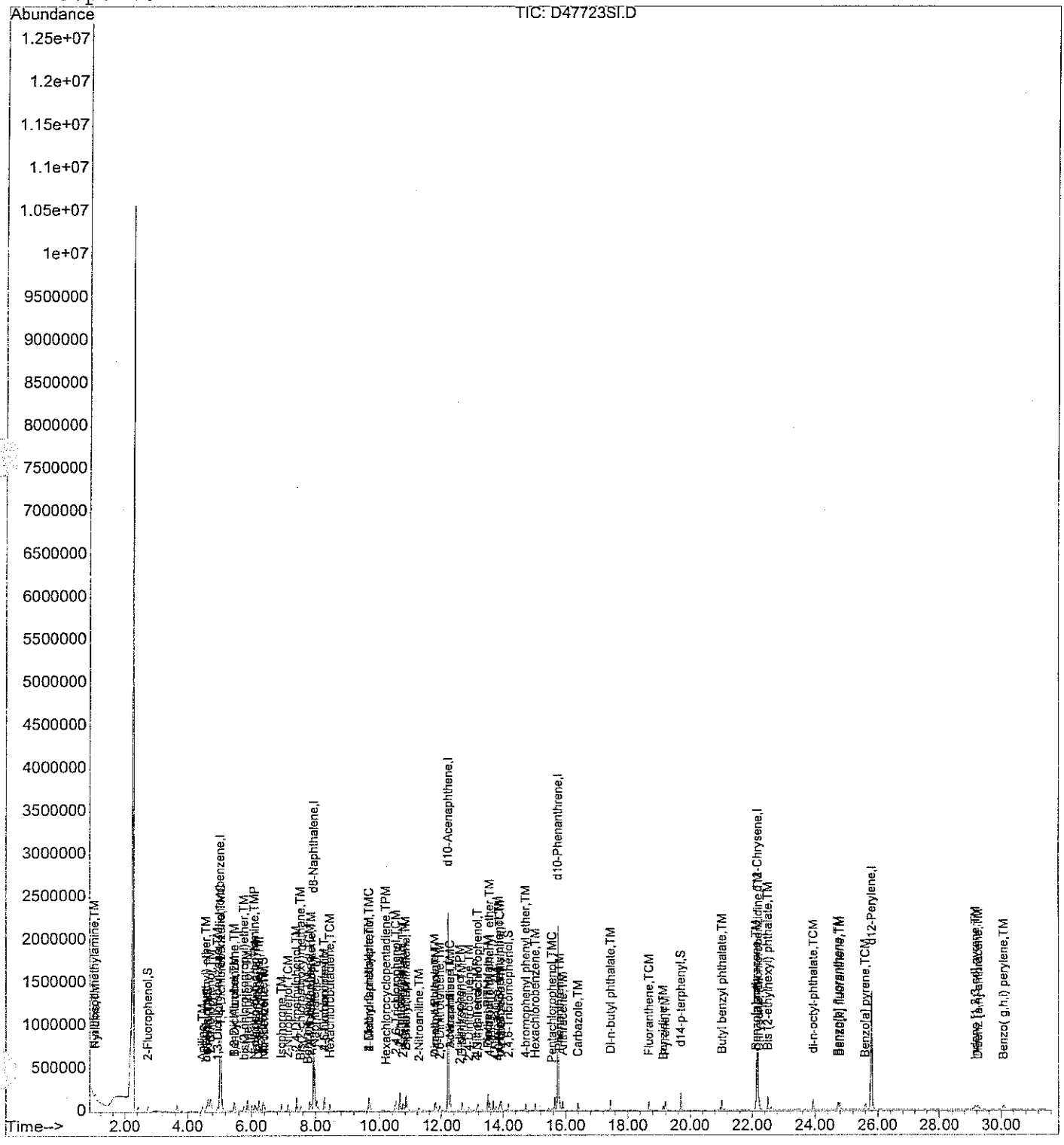
	Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
72)	Pyrene	19.19	202	83896	1.83	ug/ml	99
74)	Butyl benzyl phthalate	21.03	149	46803	1.71	ug/ml	98
75)	Benzo[a]anthracene	22.11	228	74399	1.81	ug/ml	98
76)	Chrysene	22.21	228	68382	1.77	ug/ml	99
77)	3,3'-Dichlorobenzidine	22.19	252	29168	1.98	ug/ml#	92
78)	Bis (2-ethylhexyl) phthala	22.49	149	64510	1.72	ug/ml	100
80)	di-n-octyl-phthalate	23.94	149	107973	1.69	ug/ml	97
81)	Benzo[b] fluoranthene	24.73	252	74081	1.88	ug/ml	98
82)	Benzo[k] fluoranthene	24.79	252	68826	1.82	ug/ml	97
83)	Benzo[a] pyrene	25.60	252	66524	1.78	ug/ml#	91
84)	Indeno [1,2,3-cd] pyrene	29.15	276	57599	1.52	ug/ml#	60
85)	Dibenz [a,h] anthracene	29.24	278	55419	1.52	ug/ml	97
86)	Benzo(g,h,i) perylene	30.07	276	58346	1.57	ug/ml#	84

(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : D:\HPCHEM\1\DATA\052406-D\D47723SI.D Vial: 2
Acq On : 25 May 2006 3:47 am Operator:
Sample : S6852 PT1 2/4 Inst : D Instrum
Misc : Multiplr: 1.00
MS Integration Params: RTEINT.P
Quant Time: May 26 9:23 2006 Quant Results File: S05246.RES

Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Fri May 26 09:35:43 2006
Response via : Initial Calibration

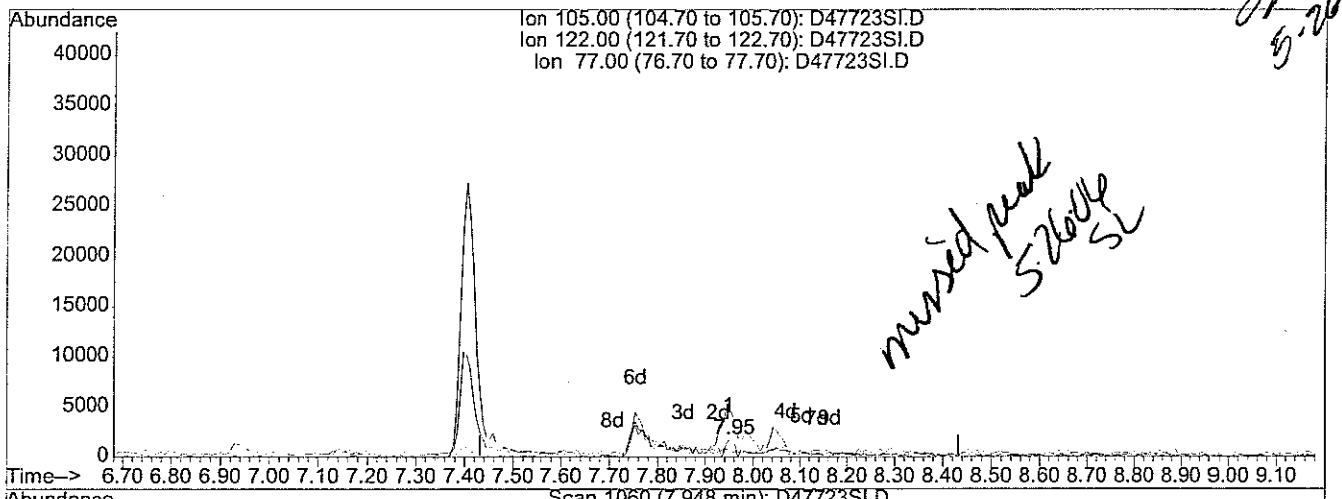


Quantitation Report (Qedit)

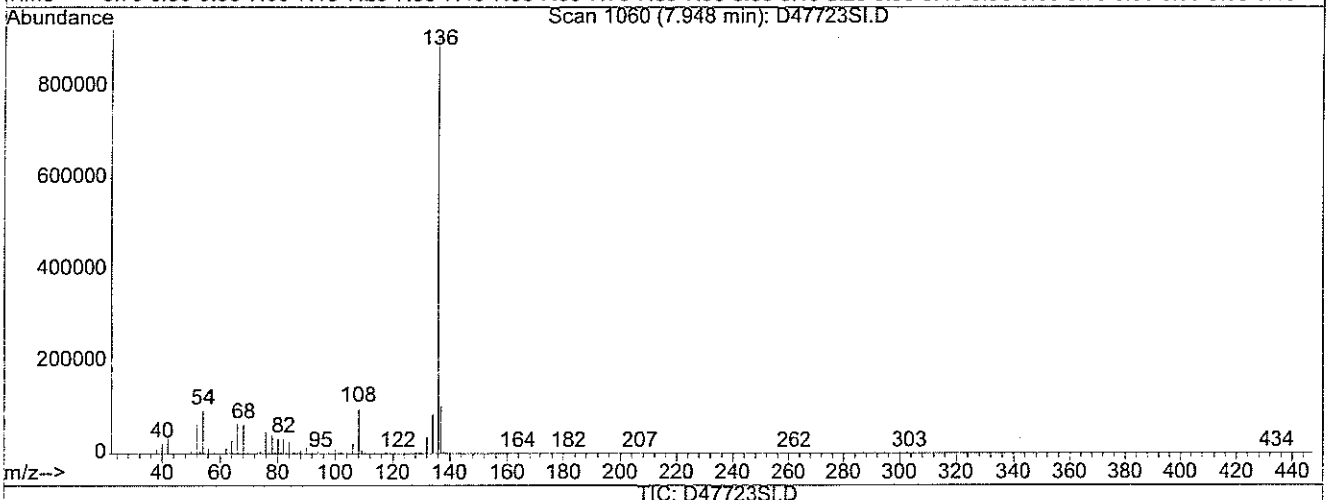
Data File : D:\HPCHEM\1\DATA\052406-D\D47723SI.D Vial: 2
Acq On : 25 May 2006 3:47 am Operator:
Sample : S6852 PT1 2/4 Inst : D Instrum
Misc : Multiplr: 1.00
MS Integration Params: RTEINT.P
Quant Time: May 26 9:22 2006 Quant Results File: temp.res

Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Fri May 26 09:21:45 2006
Response via : Multiple Level Calibration

5.26.06



missed peak 5.26.06 SL



TIC: D47723SI.D

(27) Benzoic Acid (TM)

7.95min 3.02ug/ml

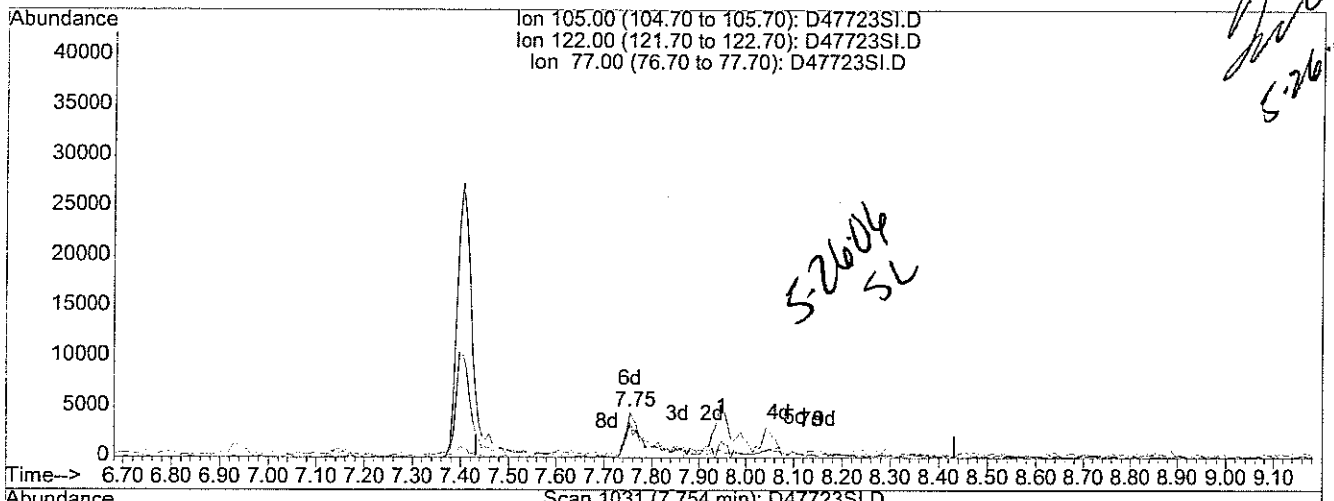
response 2193

Ion	Exp%	Act%
105.00	100	100
122.00	78.50	24.58#
77.00	70.70	457.18#
0.00	0.00	0.00

Quantitation Report (Qedit)

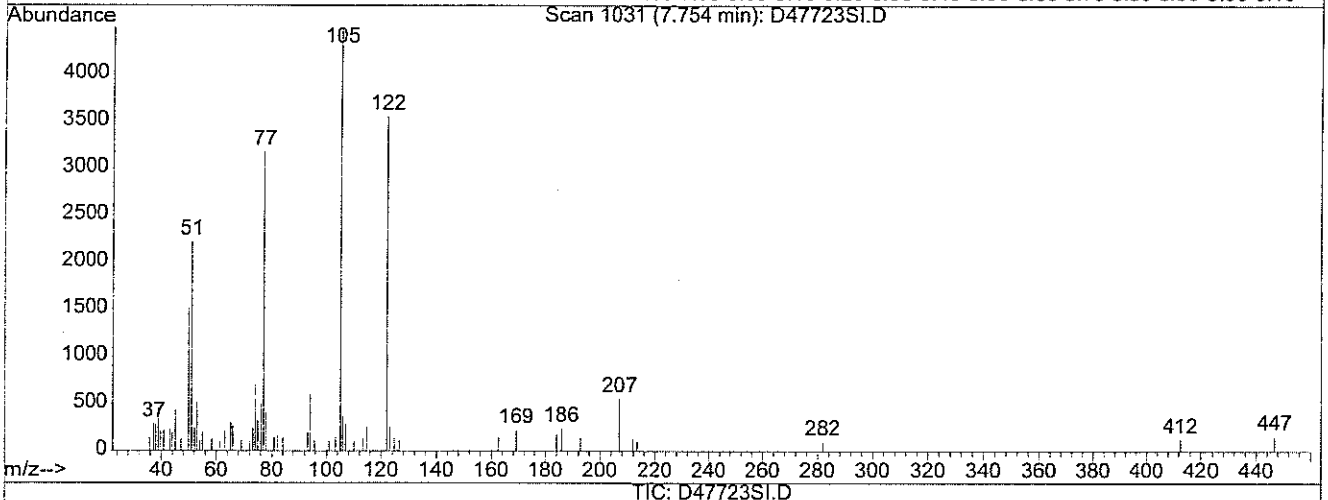
Data File : D:\HPCHEM\1\DATA\052406-D\D47723SI.D Vial: 2
Acq On : 25 May 2006 3:47 am Operator:
Sample : S6852 PT1 2/4 Inst : D Instrum
Misc : Multiplr: 1.00
MS Integration Params: RTEINT.P
Quant Time: May 26 9:22 2006 Quant Results File: temp.res

Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Fri May 26 09:21:45 2006
Response via : Multiple Level Calibration



5-26-06

5.26.06 SL



(27) Benzoic Acid (TM)

7.75min 3.97ug/ml m

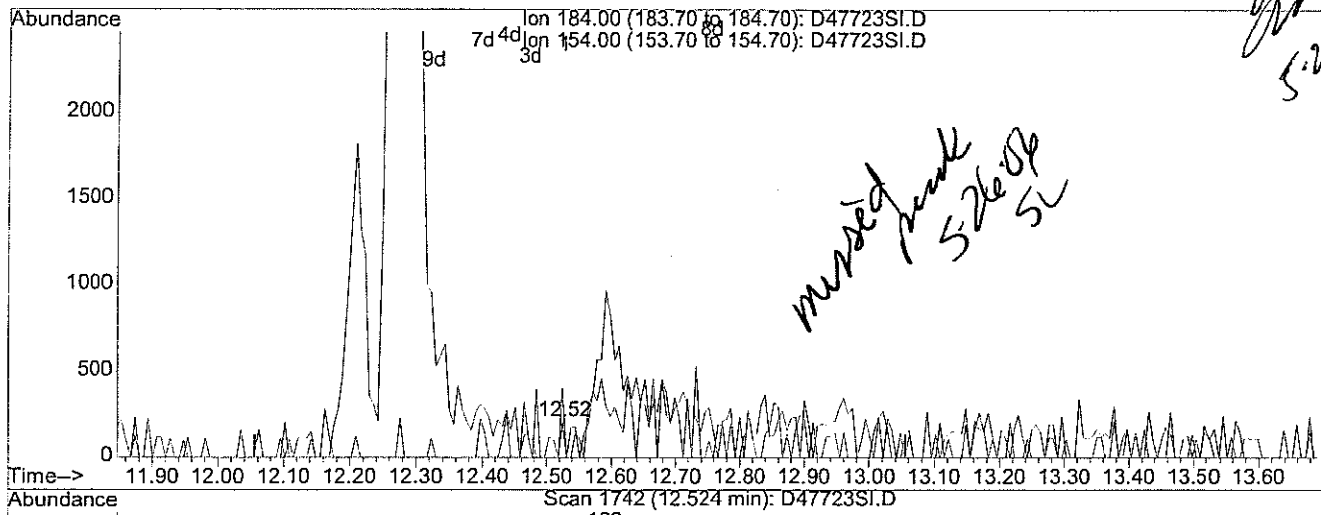
response 13546

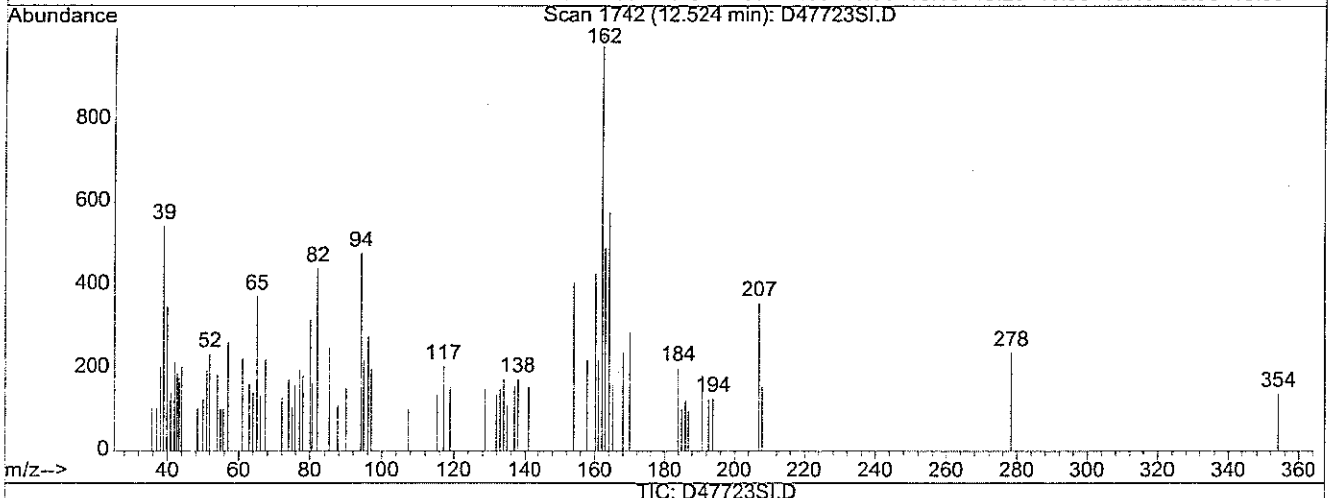
Ion	Exp%	Act%
105.00	100	100
122.00	78.50	3.98#
77.00	70.70	74.01
0.00	0.00	0.00

Quantitation Report (Qedit)

Data File : D:\HPCHEM\1\DATA\052406-D\D47723SI.D Vial: 2
Acq On : 25 May 2006 3:47 am Operator:
Sample : S6852 PT1 2/4 Inst : D Instrum
Misc : Multiplr: 1.00
MS Integration Params: RTEINT.P
Quant Time: May 26 9:22 2006 Quant Results File: temp.res

Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Fri May 26 09:21:45 2006
Response via : Multiple Level Calibration





TIC: D47723SI.D

(47) 2,4-dinitrophenol (TPM)

12.52min 4.99ug/ml

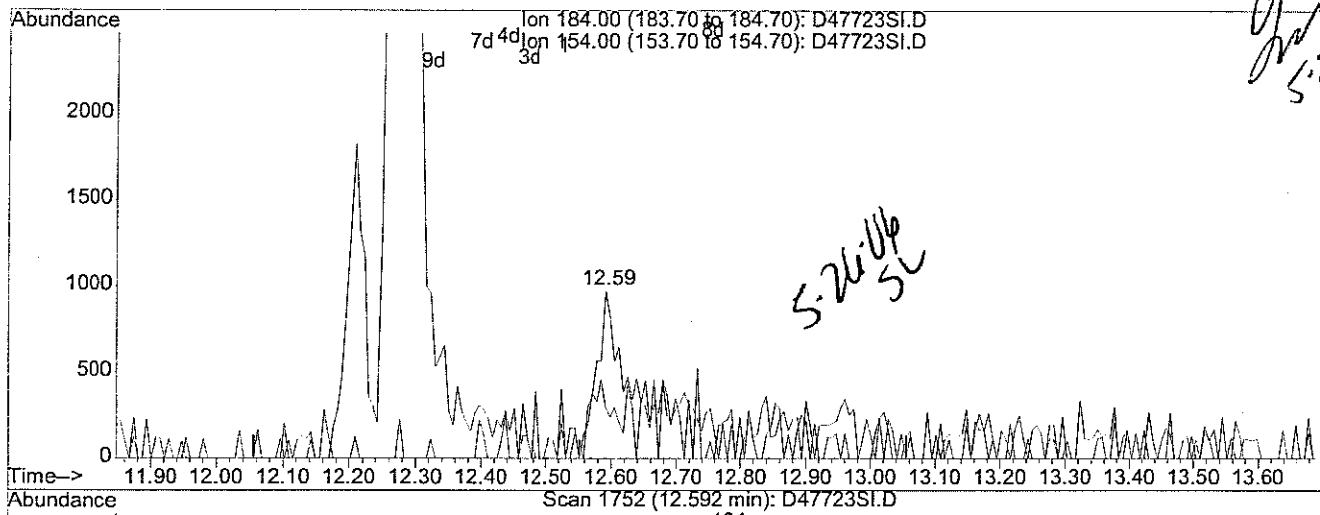
response 82

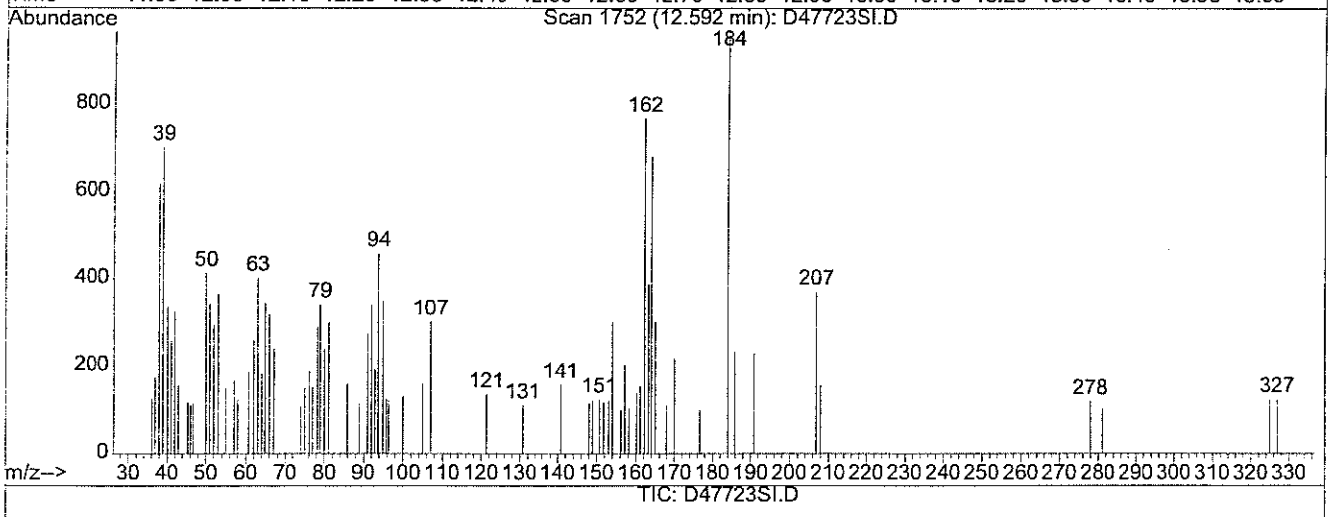
Ion	Exp%	Act%
184.00	100	100
154.00	35.30	397.56#
0.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (Qedit)

Data File : D:\HPCHEM\1\DATA\052406-D\D47723SI.D Vial: 2
Acq On : 25 May 2006 3:47 am Operator:
Sample : S6852 PT1 2/4 Inst : D Instrum
Misc : Multiplr: 1.00
MS Integration Params: RTEINT.P
Quant Time: May 26 9:23 2006 Quant Results File: temp.res

Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Fri May 26 09:21:45 2006
Response via : Multiple Level Calibration





(47) 2,4-dinitrophenol (TPM)

12.59min 5.80ug/ml m

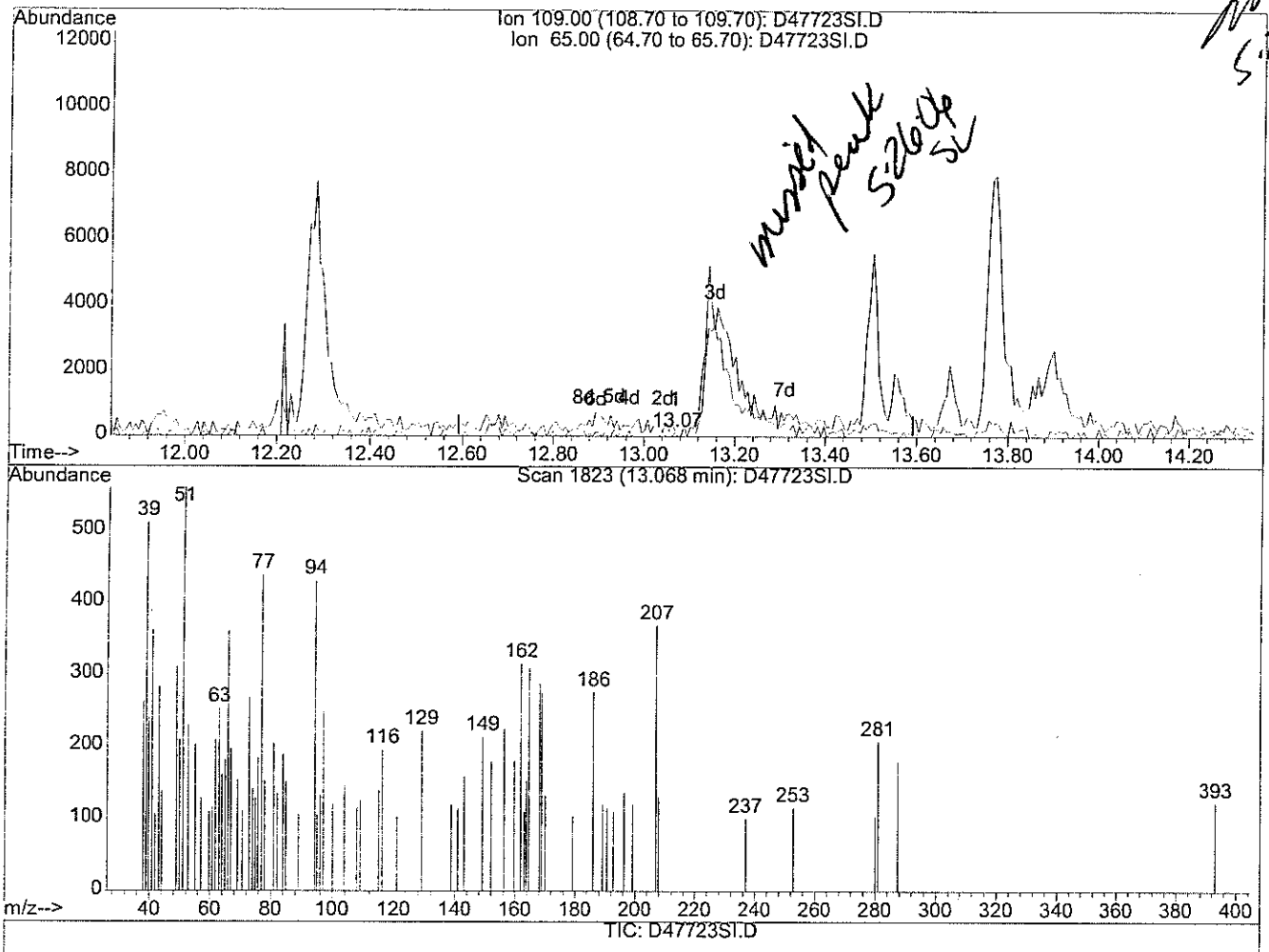
response 4691

Ion	Exp%	Act%
184.00	100	100
154.00	35.30	6.95#
0.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (Qedit)

Data File : D:\HPCHEM\1\DATA\052406-D\D47723SI.D Vial: 2
Acq On : 25 May 2006 3:47 am Operator:
Sample : S6852 PT1 2/4 Inst : D Instrum
Misc : Multiplr: 1.00
MS Integration Params: RTEINT.P
Quant Time: May 26 9:23 2006 Quant Results File: temp.res

Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Fri May 26 09:21:45 2006
Response via : Multiple Level Calibration



(49) 4-Nitrophenol (TMP)

13.07min 0.01ug/ml

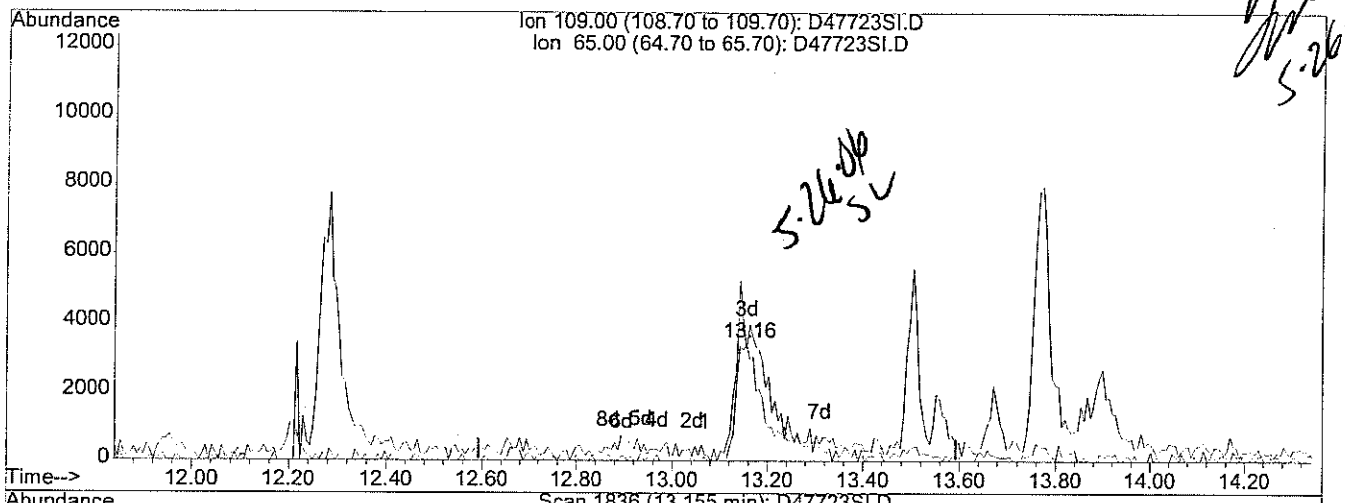
response 51

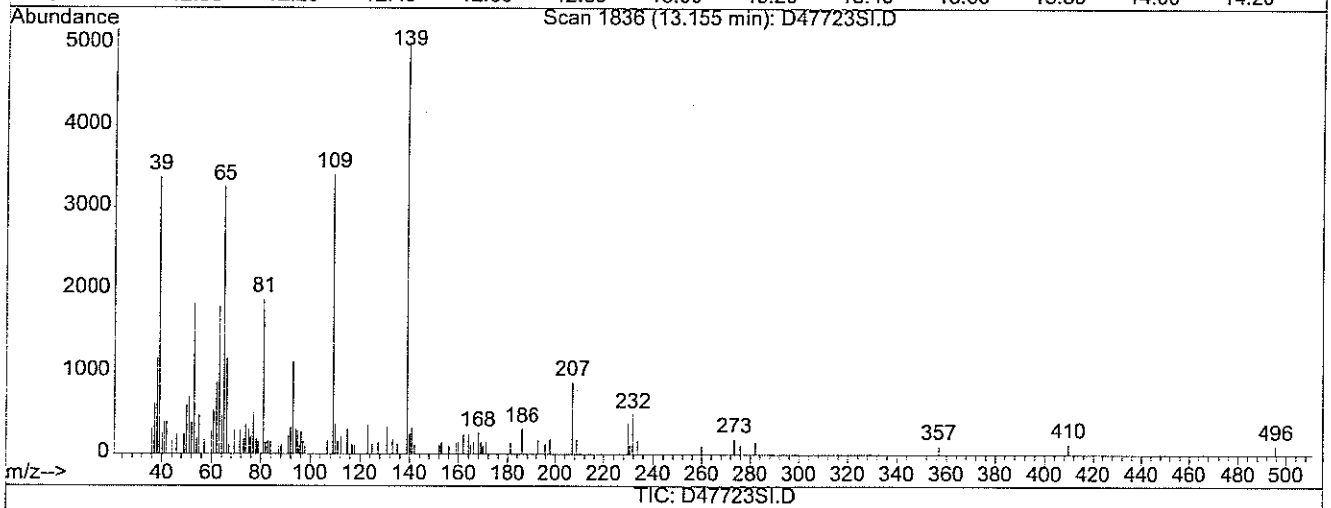
Ion	Exp%	Act%
109.00	100	100
65.00	132.70	282.35#
0.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (Qedit)

Data File : D:\HPCHEM\1\DATA\052406-D\D47723SI.D Vial: 2
Acq On : 25 May 2006 3:47 am Operator:
Sample : S6852 PT1 2/4 Inst : D Instrum
Misc : Multiplr: 1.00
MS Integration Params: RTEINT.P
Quant Time: May 26 9:23 2006 Quant Results File: temp.res

Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Fri May 26 09:21:45 2006
Response via : Multiple Level Calibration





TIC: D47723SI.D

(49) 4-Nitrophenol (TMP)

13.16min 4.04ug/ml m

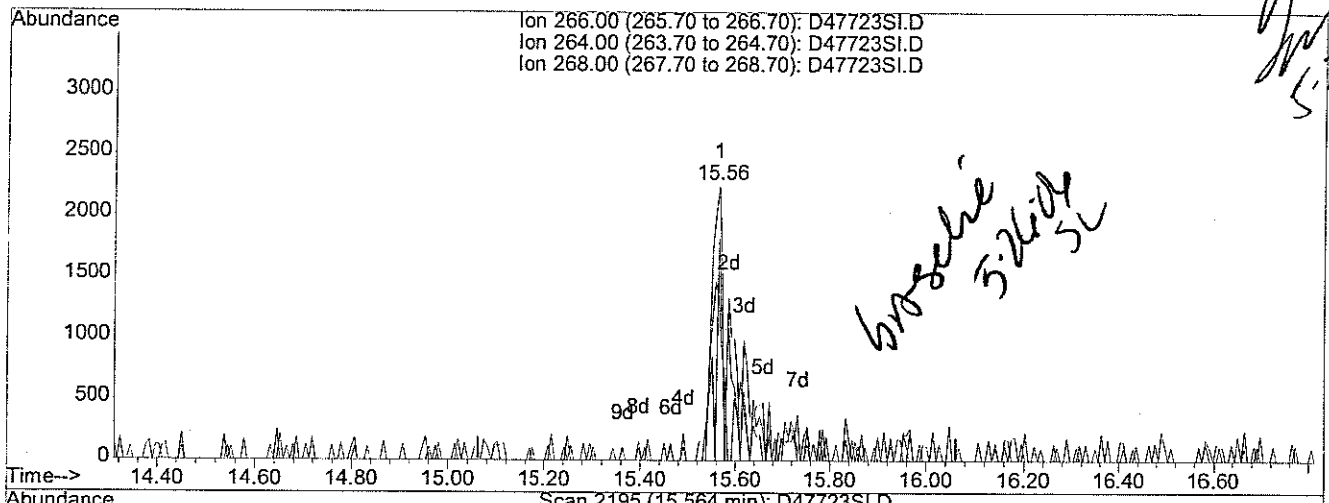
response 15483

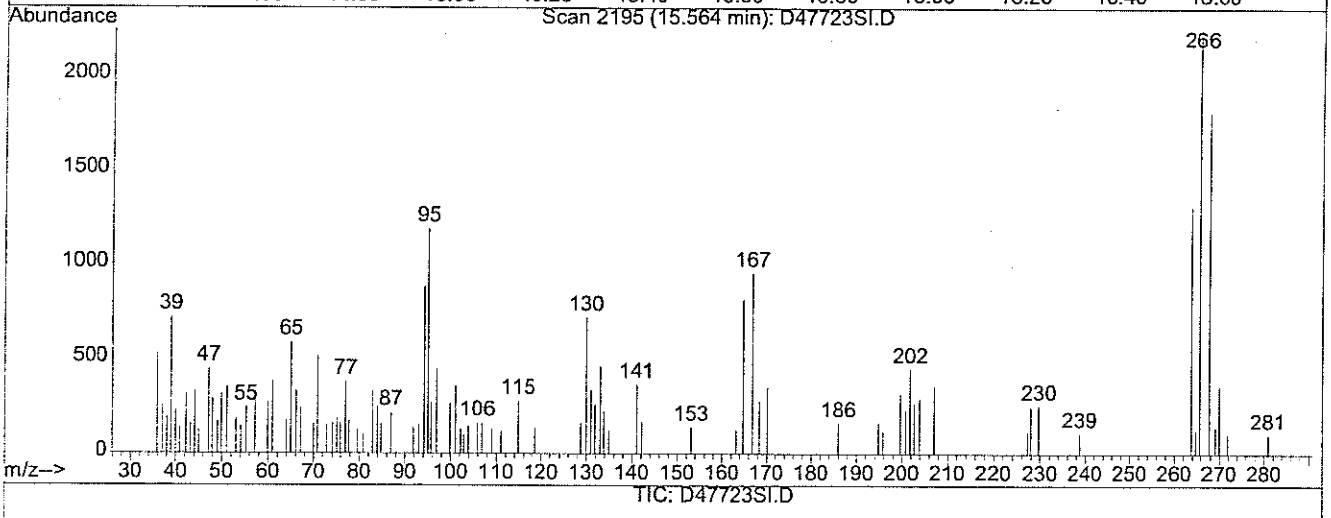
Ion	Exp%	Act%
109.00	100	100
65.00	132.70	0.93#
0.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (Qedit)

Data File : D:\HPCHEM\1\DATA\052406-D\D47723SI.D Vial: 2
Acq On : 25 May 2006 3:47 am Operator:
Sample : S6852 PT1 2/4 Inst : D Instrum
Misc : Multiplr: 1.00
MS Integration Params: RTEINT.P
Quant Time: May 26 9:23 2006 Quant Results File: temp.res

Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Fri May 26 09:21:45 2006
Response via : Multiple Level Calibration





TIC: D47723SI.D

(64) Pentachlorophenol (TMC)

15.56min 0.57ug/ml

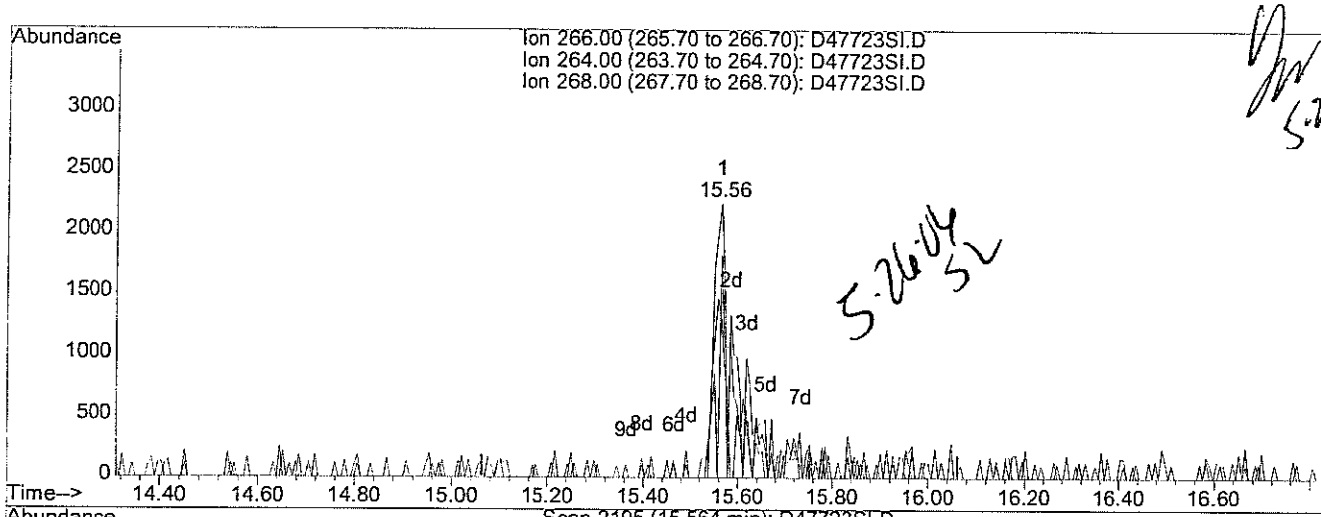
response 3301

Ion	Exp%	Act%
266.00	100	100
264.00	62.00	80.85#
268.00	69.50	101.94#
0.00	0.00	0.00

Quantitation Report (Qedit)

Data File : D:\HPCHEM\1\DATA\052406-D\D47723SI.D Vial: 2
Acq On : 25 May 2006 3:47 am Operator:
Sample : S6852 PT1 2/4 Inst : D Instrum
Misc : Multiplr: 1.00
MS Integration Params: RTEINT.P
Quant Time: May 26 9:23 2006 Quant Results File: temp.res

Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Fri May 26 09:21:45 2006
Response via : Multiple Level Calibration



(64) Pentachlorophenol (TMC)

15.56min 1.30ug/ml m

response 7480

Ion	Exp%	Act%
266.00	100	100
264.00	62.00	35.68#
268.00	69.50	44.99#
0.00	0.00	0.00

Quantitation Report (QT Reviewed)

Data File : D:\HPCHEM\1\DATA\052406-D\D47724SI.D Vial: 3
Acq On : 25 May 2006 4:20 am Operator:
Sample : S6851 PT2A 4/8 Inst : D Instrum
Misc : Multiplr: 1.00
Integration Params: RTEINT.P
Quant Time: May 26 9:24 2006 Quant Results File: S05246.RES

Quant Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Fri May 26 09:21:45 2006
Response via : Initial Calibration
DataAcq Meth : ABN

5.26.06 5.26.06

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) d4-1,4-dichlorobenzene	5.03	152	495838	40.00	ug/ml	0.00
19) d8-Naphthalene	7.95	136	1830301	40.00	ug/ml	0.00
35) d10-Acenaphthene	12.21	164	941311	40.00	ug/ml	0.00
61) d10-Phenanthrene	15.75	188	1609031	40.00	ug/ml	0.00
71) d12-Chrysene	22.16	240	1433198	40.00	ug/ml	0.00
79) d12-Perylene	25.80	264	1362420	40.00	ug/ml	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	2.73	112	150717	8.50	ug/ml	0.00
Spiked Amount 100.000	Range 28 - 110		Recovery = 8.50%#			
5) d5-Phenol	4.62	99	175559	8.37	ug/ml	0.01
Spiked Amount 100.000	Range 22 - 110		Recovery = 8.37%#			
20) d5-nitrobenzene	6.36	82	154035	7.42	ug/ml	-0.01
Spiked Amount 50.000	Range 45 - 107		Recovery = 14.84%#			
39) 2-Fluorobiphenyl	10.64	172	245118	7.87	ug/ml	0.00
Spiked Amount 50.000	Range 44 - 102		Recovery = 15.74%#			
60) 2,4,6-Tribromophenol	14.17	330	37363	6.16	ug/ml	0.00
Spiked Amount 100.000	Range 63 - 127		Recovery = 6.16%#			
73) d14-p-terphenyl	19.70	244	236213	7.65	ug/ml	-0.01
Spiked Amount 50.000	Range 33 - 132		Recovery = 15.30%#			
Target Compounds						Qvalue
2) Pyridine	1.02	79	71406	2.65	ug/ml#	100
3) N-nitrosodimethylamine	1.00	42	57877	3.72	ug/ml	93
6) Aniline	4.46	93	118452	4.06	ug/ml	97
7) Phenol	4.64	94	195480	8.57	ug/ml	97
8) 2-Chlorophenol	4.74	128	146867	7.79	ug/ml	99
9) bis(2-chloroethyl) ether	4.58	93	73885	3.85	ug/ml	83
10) 1,3-Dichlorobenzene	4.95	146	74657	3.78	ug/ml	97
11) 1,4-Dichlorobenzene	5.07	146	74331	3.70	ug/ml#	83
12) 1,2-Dichlorobenzene	5.48	146	70071	3.71	ug/ml	95
13) Benzyl Alcohol	5.46	108	49568	4.45	ug/ml	92
14) 2-Methylphenol	5.88	108	138457	8.82	ug/ml	99
15) bis(2-chloroisopropyl)ethe	5.78	45	123053	3.31	ug/ml	98
16) 3+4-Methylphenol	6.23	108	143186	8.87	ug/ml	99
17) Hexachloroethane	6.13	117	34996	4.22	ug/ml	99
18) N-nitroso-di-n-propylamine	6.10	70	48216	3.18	ug/ml	97
21) Nitrobenzene	6.40	123	39837	4.05	ug/ml	92
22) Isophorone	6.93	82	131782	3.67	ug/ml	99
23) 2-Nitrophenol	7.15	139	78715	7.22	ug/ml	96
24) 2,4-Dimethylphenol	7.40	107	148704	9.23	ug/ml	98
25) Bis(2-chloroethoxy)methane	7.54	93	86566	4.04	ug/ml	99

(#) = qualifier out of range (m) = manual integration

Quantitation Report (QT Reviewed)

Data File : D:\HPCHEM\1\DATA\052406-D\D47724SI.D Vial: 3
Acq On : 25 May 2006 4:20 am Operator:
Sample : S6851 PT2A 4/8 Inst : D Instrum
Misc : Multiplr: 1.00
Integration Params: RTEINT.P
Quant Time: May 26 9:24 2006 Quant Results File: S05246.RES

Quant Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Fri May 26 09:21:45 2006
Response via : Initial Calibration
DataAcq Meth : ABN

Compound	R.T.	QIon	Response	Conc Unit	Qvalue
26) 2,4-Dichlorophenol	7.81	162	101574	8.03 ug/ml	98
27) Benzoic Acid	7.78	105	47059m	6.91 ug/ml	
28) 1,2,4-Trichlorobenzene	7.87	180	55430	4.07 ug/ml	99
29) Naphthalene	7.99	128	197102	4.01 ug/ml	100
30) 2,6-Dichlorophenol	8.28	162	101420	7.98 ug/ml	100
31) 4-Chloroaniline	8.27	127	82690	3.79 ug/ml	96
32) Hexachlorobutadiene	8.45	225	30362	3.81 ug/ml	98
33) 4-Chloro-3-methylphenol	9.66	107	118769	8.92 ug/ml	99
34) 2-Methylnaphthalene	9.65	141	109962	3.95 ug/ml	96
36) Hexachlorocyclopentadiene	10.21	237	1497	4.21 ug/ml#	72
37) 2,4,6-Trichlorophenol	10.52	196	64549	6.61 ug/ml	97
38) 2,4,5-Trichlorophenol	10.72	196	66722	6.63 ug/ml	95
40) 2-Chloronaphthalene	10.82	162	108744	3.77 ug/ml	99
41) 2-Nitroaniline	11.25	138	45133	3.27 ug/ml	95
42) 2,6-Dinitrotoluene	11.95	165	30007	3.47 ug/ml	97
43) Dimethyl Phthalate	11.78	163	129470	3.79 ug/ml	99
44) Acenaphthylene	11.82	152	171234	3.56 ug/ml	97
45) 3-Nitroaniline	12.27	138	36590	3.36 ug/ml	95
46) Acenaphthene	12.27	154	94924	3.55 ug/ml	97
47) 2,4-dinitrophenol	12.58	184	15516	7.69 ug/ml#	65
48) Dibenzofuran	12.67	168	157980	3.90 ug/ml	97
49) 4-Nitrophenol	13.13	109	32311	8.42 ug/ml	97
50) 2,4-Dinitrotoluene	12.89	165	41033	3.71 ug/ml	96
51) 2,3,4,6-Tetrachlorophenol	13.18	232	41879	5.54 ug/ml	98
52) Fluorene	13.49	166	120488	3.87 ug/ml	100
53) Diethyl Phthalate	13.51	149	134749	3.74 ug/ml	96
54) 4-Chlorophenyl phenyl eth	13.56	204	54757	3.50 ug/ml#	88
55) 4-Nitroaniline	13.77	138	37993	3.39 ug/ml	99
56) 4,6-Dinitro-2-methylphenol	13.86	198	37828	7.49 ug/ml	100
57) n-nitrosodiphenylamine	13.90	169	81301	3.85 ug/ml	98
58) Azobenzene	13.93	77	179584	3.65 ug/ml	98
59) Biphenyl	10.83	154	136967	3.63 ug/ml	99
62) 4-bromophenyl phenyl ether	14.72	248	34113	3.50 ug/ml	96
63) Hexachlorobenzene	15.02	284	39749	3.59 ug/ml	96
64) Pentachlorophenol	15.56	266	21376m	3.67 ug/ml	
65) Phenanthrene	15.79	178	159688	3.77 ug/ml	99
66) Anthracene	15.90	178	162394	3.80 ug/ml	99
67) Carbazole	16.39	167	160279	3.90 ug/ml	100
68) Di-n-butyl phthalate	17.45	149	220918	3.59 ug/ml	99
69) Fluoranthene	18.67	202	163233	3.57 ug/ml	99
70) Benzidine	19.12	184	94483	4.90 ug/ml	98

(#) = qualifier out of range (m) = manual integration

Quantitation Report (QT Reviewed)

Data File : D:\HPCHEM\1\DATA\052406-D\D47724SI.D Vial: 3
Acq On : 25 May 2006 4:20 am Operator:
Sample : S6851 PT2A 4/8 Inst : D Instrum
Misc : Multiplr: 1.00
Integration Params: RTEINT.P
Quant Time: May 26 9:24 2006 Quant Results File: S05246.RES

Quant Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Fri May 26 09:21:45 2006
Response via : Initial Calibration
DataAcq Meth : ABN

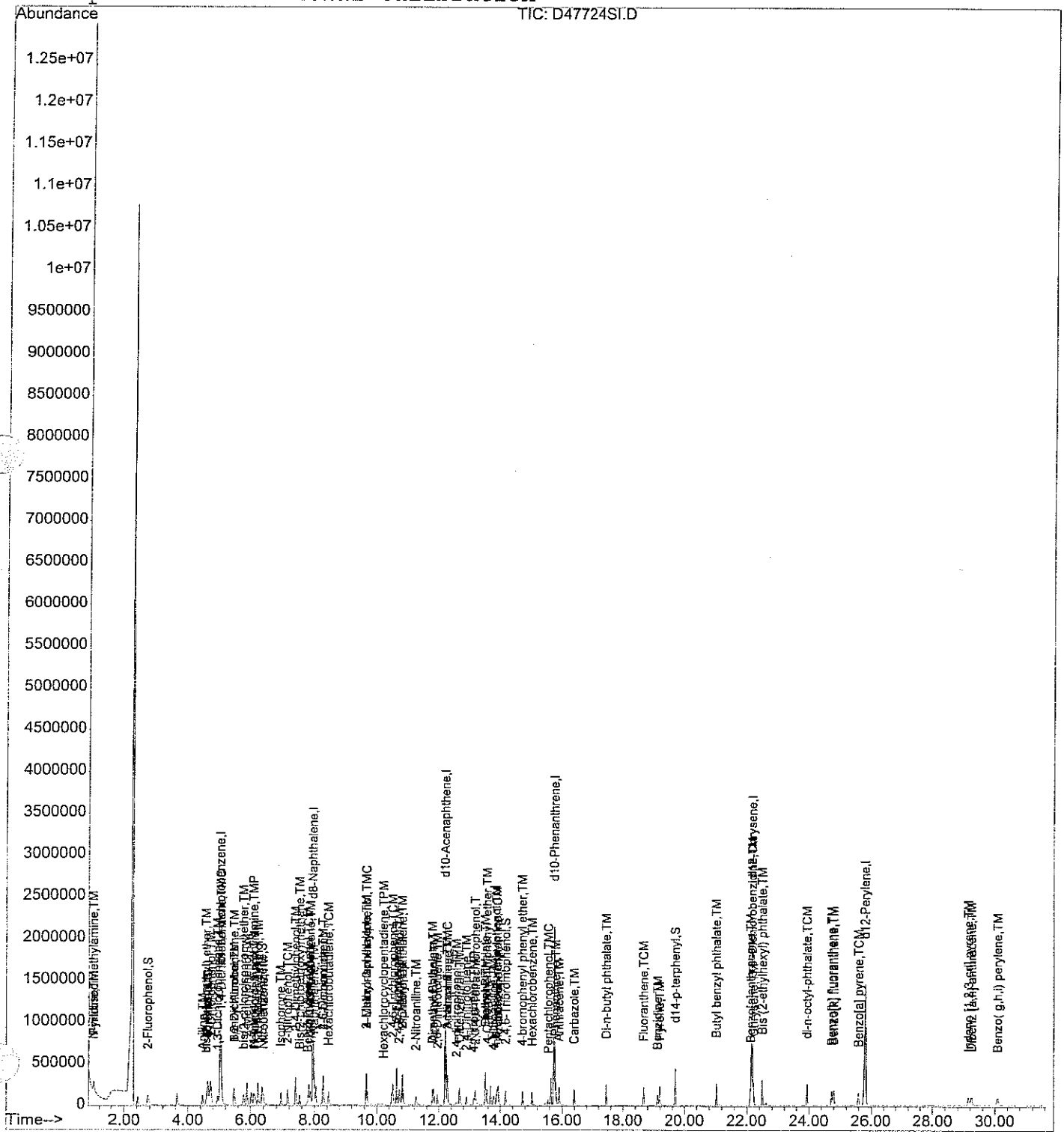
Compound	R.T.	QIon	Response	Conc Unit	Qvalue
72) Pyrene	19.19	202	168806	3.60 ug/ml	99
74) Butyl benzyl phthalate	21.02	149	101105	3.61 ug/ml	96
75) Benzo[a]anthracene	22.11	228	151822	3.62 ug/ml	99
76) Chrysene	22.21	228	141080	3.57 ug/ml	99
77) 3,3'-Dichlorobenzidine	22.18	252	62086	4.12 ug/ml#	95
78) Bis (2-ethylhexyl) phthala	22.49	149	136204	3.56 ug/ml	99
80) di-n-octyl-phthalate	23.94	149	232005	3.59 ug/ml	100
81) Benzo[b] fluoranthene	24.73	252	142919	3.59 ug/ml	98
82) Benzo[k] fluoranthene	24.79	252	143180	3.76 ug/ml	98
83) Benzo[a] pyrene	25.60	252	140931	3.73 ug/ml#	91
84) Indeno [1,2,3-cd] pyrene	29.15	276	125069	3.27 ug/ml#	64
85) Dibenz [a,h] anthracene	29.24	278	120686	3.29 ug/ml	97
86) Benzo(g,h,i) perylene	30.08	276	127796	3.40 ug/ml#	85

(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : D:\HPCHEM\1\DATA\052406-D\D47724SI.D Vial: 3
Acq On : 25 May 2006 4:20 am Operator:
Sample : S6851 PT2A 4/8 Inst : D Instrum
Misc : Multiplr: 1.00
MS Integration Params: RTEINT.P
Quant Time: May 26 9:24 2006 Quant Results File: S05246.RES

Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Fri May 26 09:35:43 2006
Response via : Initial Calibration

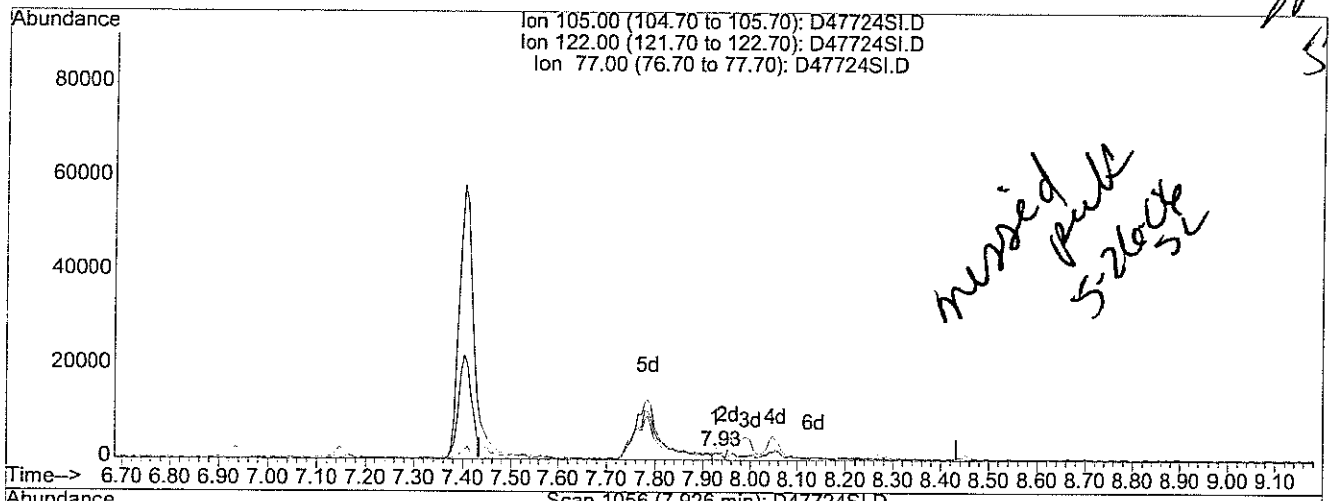


Quantitation Report (Qedit)

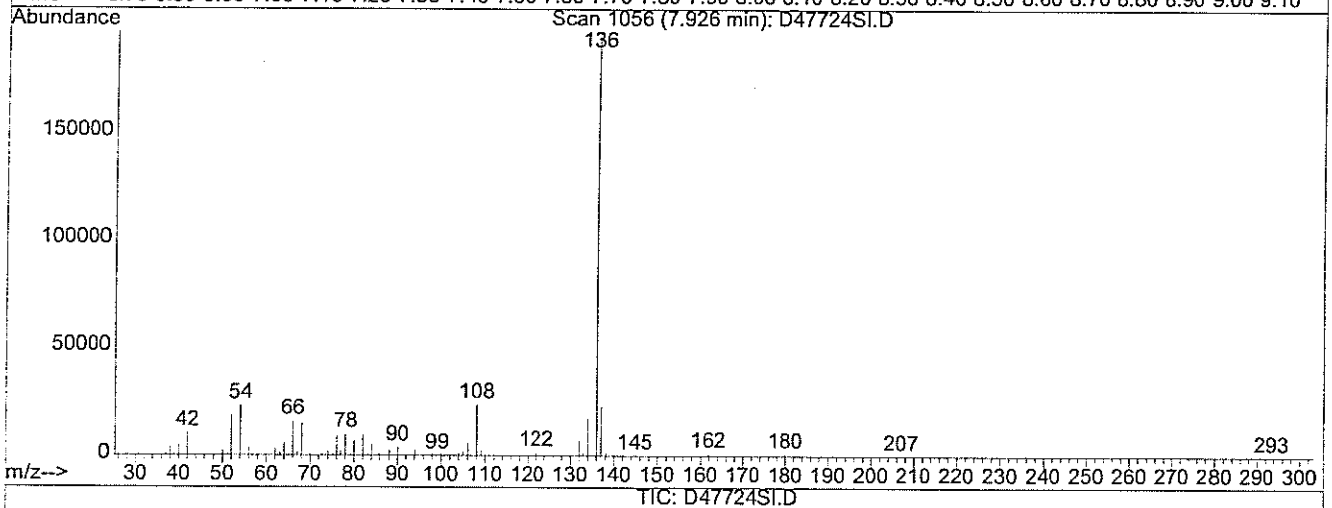
Data File : D:\HPCHEM\1\DATA\052406-D\D47724SI.D Vial: 3
Acq On : 25 May 2006 4:20 am Operator:
Sample : S6851 PT2A 4/8 Inst : D Instrum
Misc : Multiplr: 1.00
MS Integration Params: RTEINT.P
Quant Time: May 26 9:22 2006 Quant Results File: temp.res

Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Fri May 26 09:21:45 2006
Response via : Multiple Level Calibration

5-26-06



missed peak 5-26-06 SL



TIC: D47724SI.D

(27) Benzoic Acid (TM)

7.93min 2.99ug/ml

response 1850

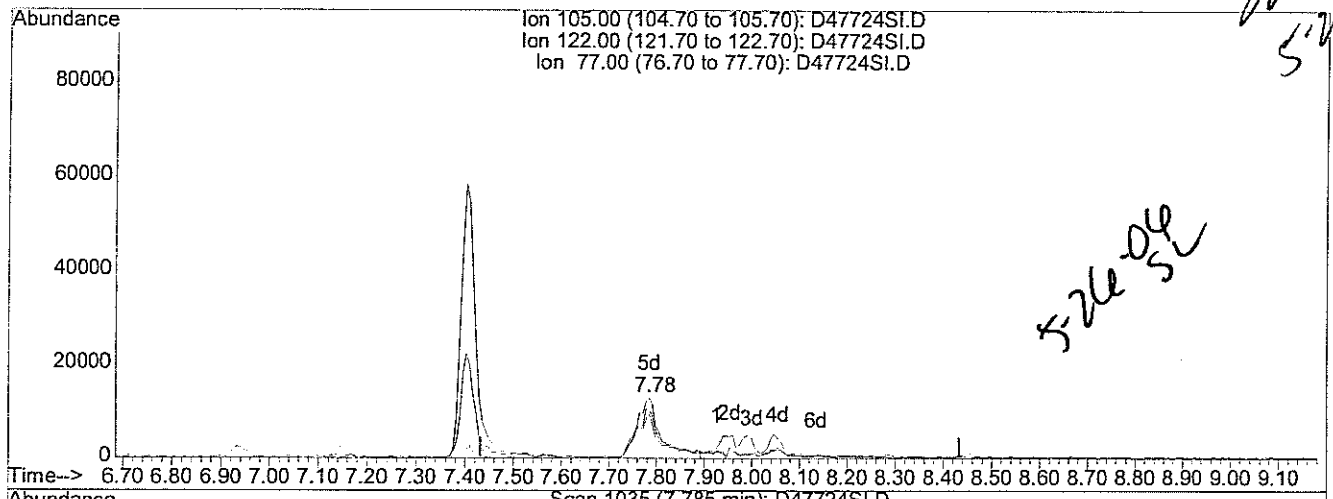
Ion	Exp%	Act%
105.00	100	100
122.00	78.50	0.00#
77.00	70.70	0.00#
0.00	0.00	0.00

Quantitation Report (Qedit)

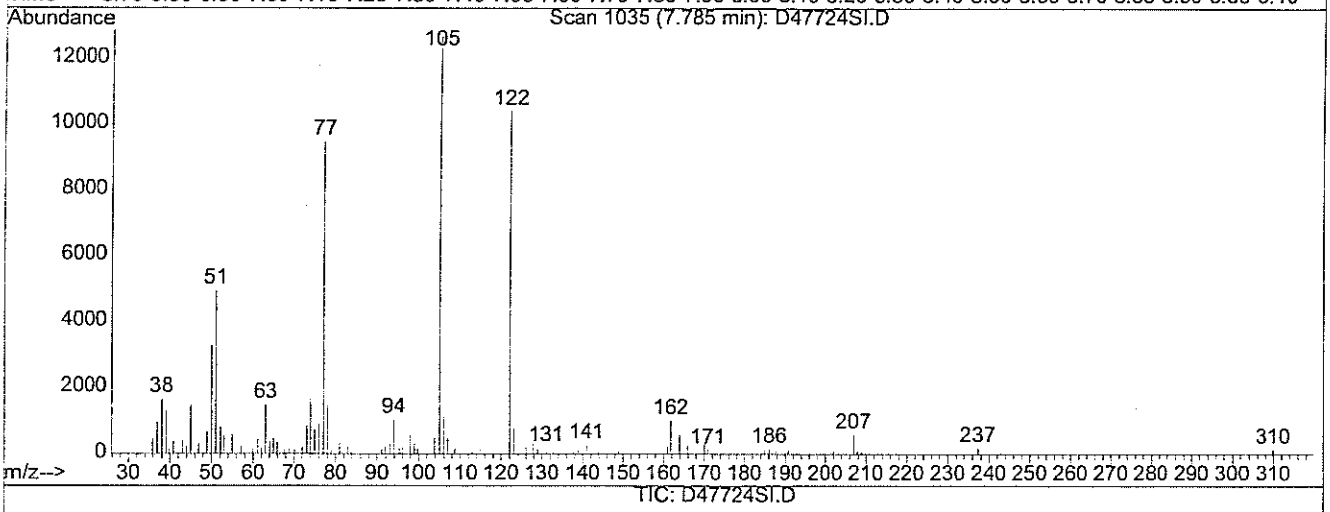
Data File : D:\HPCHEM\1\DATA\052406-D\D47724SI.D Vial: 3
Acq On : 25 May 2006 4:20 am Operator:
Sample : S6851 PT2A 4/8 Inst : D Instrum
Misc : Multiplr: 1.00
MS Integration Params: RTEINT.P
Quant Time: May 26 9:24 2006 Quant Results File: temp.res

Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Fri May 26 09:21:45 2006
Response via : Multiple Level Calibration

5-26-06



5-26-06 SV



TIC: D47724SI.D

(27) Benzoic Acid (TM)

7.78min 6.91ug/ml m

response 47059

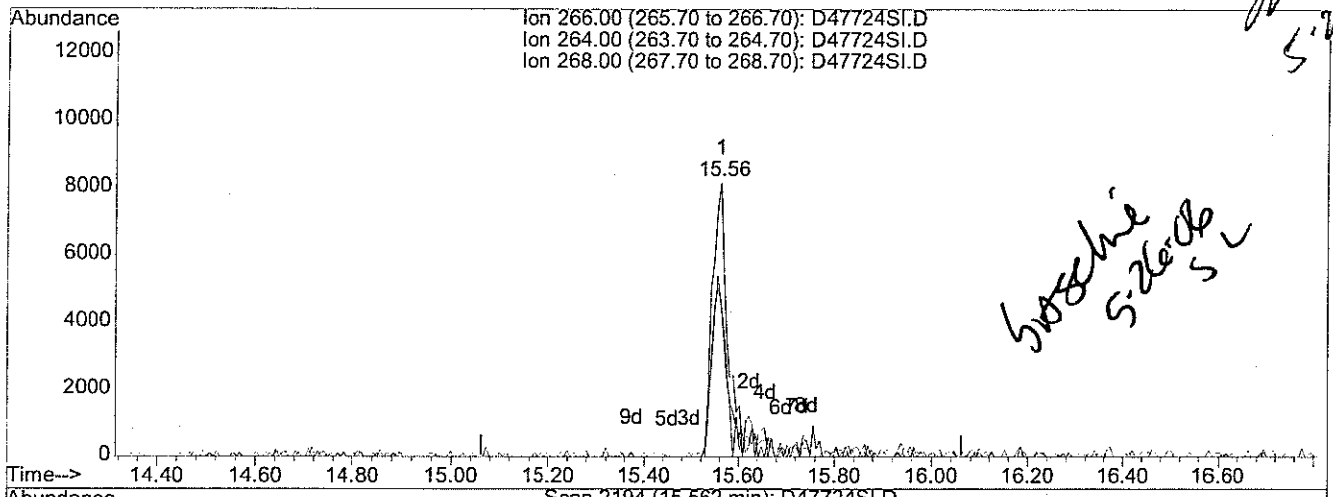
Ion	Exp%	Act%
105.00	100	100
122.00	78.50	0.00#
77.00	70.70	0.00#
0.00	0.00	0.00

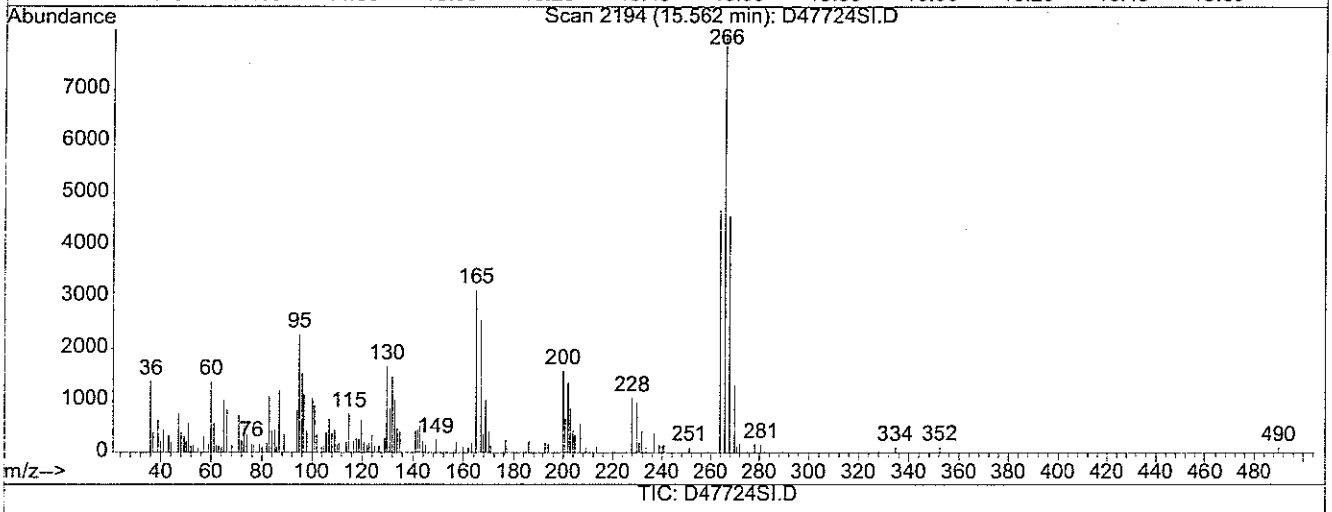
Quantitation Report (Qedit)

Data File : D:\HPCHEM\1\DATA\052406-D\D47724SI.D Vial: 3
Acq On : 25 May 2006 4:20 am Operator:
Sample : S6851 PT2A 4/8 Inst : D Instrum
Misc : Multiplr: 1.00
MS Integration Params: RTEINT.P
Quant Time: May 26 9:24 2006 Quant Results File: temp.res

Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Fri May 26 09:21:45 2006
Response via : Multiple Level Calibration

5-26-06





TIC: D47724SI.D

(64) Pentachlorophenol (TMC)

15.56min 3.10ug/ml

response 18076

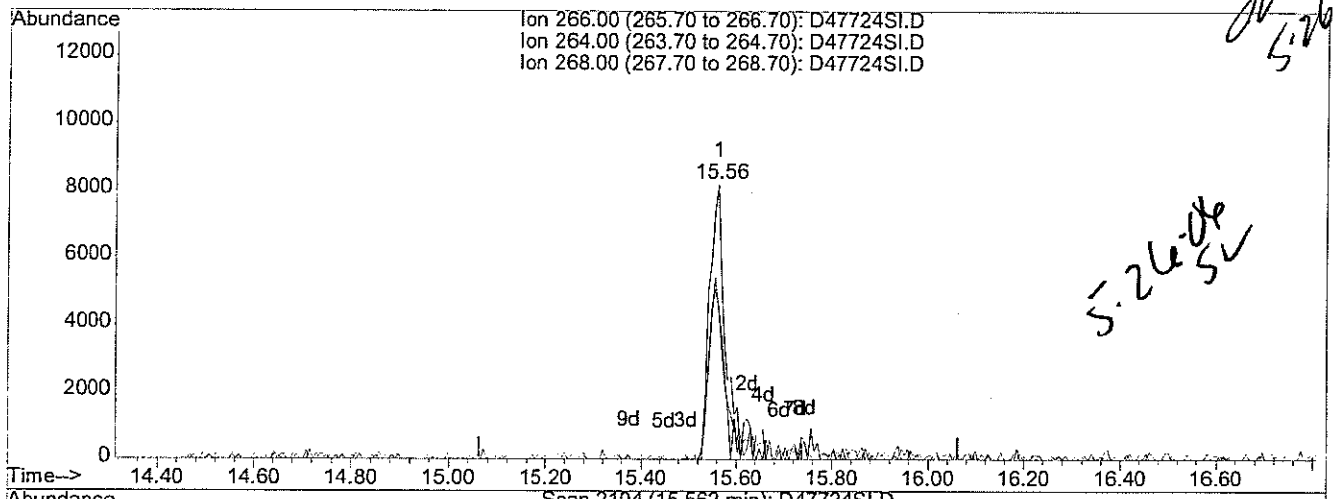
Ion	Exp%	Act%
266.00	100	100
264.00	62.00	63.48
268.00	69.50	59.61
0.00	0.00	0.00

Quantitation Report (Qedit)

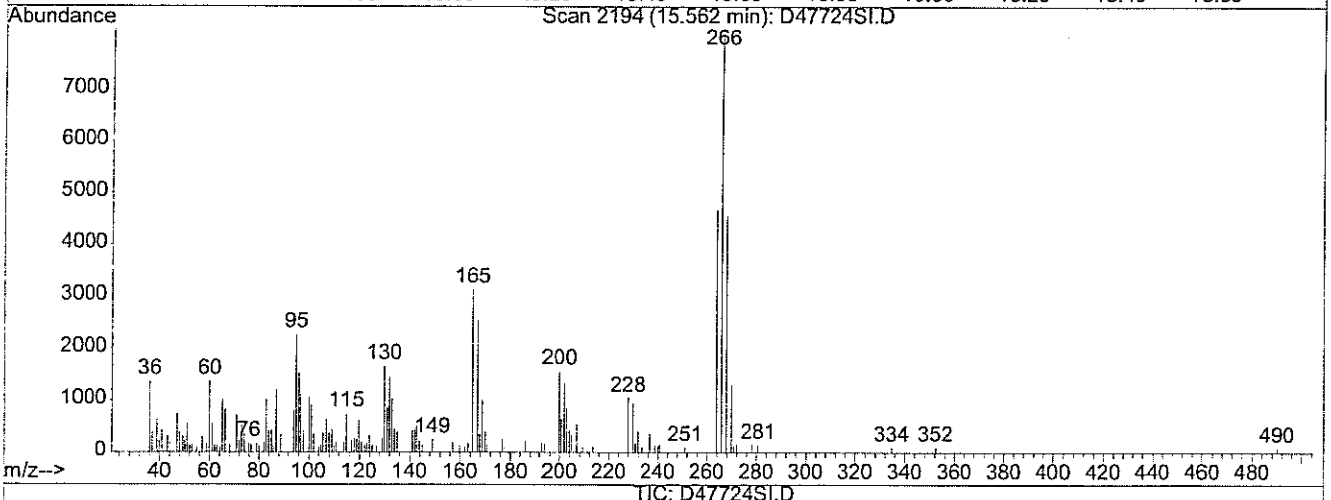
Data File : D:\HPCHEM\1\DATA\052406-D\D47724SI.D Vial: 3
Acq On : 25 May 2006 4:20 am Operator:
Sample : S6851 PT2A 4/8 Inst : D Instrum
Misc : Multiplr: 1.00
MS Integration Params: RTEINT.P
Quant Time: May 26 9:24 2006 Quant Results File: temp.res

Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Fri May 26 09:21:45 2006
Response via : Multiple Level Calibration

5.26.06



5.26.06 SV



TIC: D47724SI.D

(64) Pentachlorophenol (TMC)

15.56min 3.67ug/ml m

response 21376

Ion	Exp%	Act%
266.00	100	100
264.00	62.00	53.68
268.00	69.50	50.41#
0.00	0.00	0.00

Quantitation Report (QT Reviewed)

Data File : D:\HPCHEM\1\DATA\052406-D\D47725SI.D Vial: 4
Acq On : 25 May 2006 4:53 am Operator:
Sample : S6850 PT2 10/20 Inst : D Instrum
Misc : Multiplr: 1.00
Integration Params: RTEINT.P
Quant Time: May 26 9:22 2006 Quant Results File: S05246.RES

Quant Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Fri May 26 09:21:45 2006
Response via : Initial Calibration
DataAcq Meth : ABN

5.26.06 5.26.06 SL

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) d4-1,4-dichlorobenzene	5.04	152	492927	40.00	ug/ml	0.00
19) d8-Naphthalene	7.95	136	1801957	40.00	ug/ml	0.00
35) d10-Acenaphthene	12.21	164	930235	40.00	ug/ml	0.00
61) d10-Phenanthrene	15.75	188	1577636	40.00	ug/ml	0.00
71) d12-Chrysene	22.17	240	1404558	40.00	ug/ml	0.00
79) d12-Perylene	25.80	264	1346263	40.00	ug/ml	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	2.73	112	374538	21.25	ug/ml	0.00
Spiked Amount 100.000	Range 28 - 110		Recovery = 21.25%#			
5) d5-Phenol	4.61	99	443793	21.28	ug/ml	0.00
Spiked Amount 100.000	Range 22 - 110		Recovery = 21.28%#			
20) d5-nitrobenzene	6.36	82	405796	19.86	ug/ml	0.00
Spiked Amount 50.000	Range 45 - 107		Recovery = 39.72%#			
39) 2-Fluorobiphenyl	10.64	172	613860	19.96	ug/ml	0.00
Spiked Amount 50.000	Range 44 - 102		Recovery = 39.92%#			
60) 2,4,6-Tribromophenol	14.18	330	95463	15.93	ug/ml	0.00
Spiked Amount 100.000	Range 63 - 127		Recovery = 15.93%#			
73) d14-p-terphenyl	19.70	244	585201	19.34	ug/ml	-0.01
Spiked Amount 50.000	Range 33 - 132		Recovery = 38.68%			
Target Compounds						Qvalue
2) Pyridine	1.01	79	178242	6.65	ug/ml#	100
3) N-nitrosodimethylamine	1.00	42	144008	9.30	ug/ml	98
6) Aniline	4.47	93	296676	10.23	ug/ml	98
7) Phenol	4.63	94	477169	21.03	ug/ml	100
8) 2-Chlorophenol	4.73	128	372453	19.87	ug/ml	98
9) bis(2-chloroethyl) ether	4.59	93	201410	10.57	ug/ml	99
10) 1,3-Dichlorobenzene	4.95	146	178609	9.09	ug/ml	98
11) 1,4-Dichlorobenzene	5.07	146	186487	9.34	ug/ml	94
12) 1,2-Dichlorobenzene	5.48	146	175017	9.32	ug/ml	97
13) Benzyl Alcohol	5.46	108	125978	11.38	ug/ml	93
14) 2-Methylphenol	5.87	108	351010	22.50	ug/ml	99
15) bis(2-chloroisopropyl)ethe	5.77	45	305745	8.27	ug/ml	100
16) 3+4-Methylphenol	6.23	108	356922	22.25	ug/ml	98
17) Hexachloroethane	6.14	117	87539	10.61	ug/ml	99
18) N-nitroso-di-n-propylamine	6.10	70	116842	7.75	ug/ml	95
21) Nitrobenzene	6.40	123	101284	10.45	ug/ml	98
22) Isophorone	6.94	82	332184	9.39	ug/ml	99
23) 2-Nitrophenol	7.14	139	201941	18.81	ug/ml	96
24) 2,4-Dimethylphenol	7.40	107	376607	23.75	ug/ml	97
25) Bis(2-chloroethoxy)methane	7.54	93	225871	10.71	ug/ml	98

(#) = qualifier out of range (m) = manual integration

Quantitation Report (QT Reviewed)

Data File : D:\HPCHEM\1\DATA\052406-D\D47725SI.D Vial: 4
Acq On : 25 May 2006 4:53 am Operator:
Sample : S6850 PT2 10/20 Inst : D Instrum
Misc : Multiplr: 1.00
Integration Params: RTEINT.P
Quant Time: May 26 9:22 2006 Quant Results File: S05246.RES

Quant Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Fri May 26 09:21:45 2006
Response via : Initial Calibration
DataAcq Meth : ABN

Compound	R.T.	QIon	Response	Conc Unit	Qvalue
26) 2,4-Dichlorophenol	7.81	162	257051	20.63 ug/ml	99
27) Benzoic Acid	7.85	105	168098	17.55 ug/ml	98
28) 1,2,4-Trichlorobenzene	7.87	180	135701	10.11 ug/ml	100
29) Naphthalene	8.00	128	494448	10.22 ug/ml	99
30) 2,6-Dichlorophenol	8.28	162	258096	20.62 ug/ml	100
31) 4-Chloroaniline	8.27	127	215303	10.02 ug/ml	97
32) Hexachlorobutadiene	8.45	225	74024	9.45 ug/ml	99
33) 4-Chloro-3-methylphenol	9.66	107	299167	22.83 ug/ml	99
34) 2-Methylnaphthalene	9.65	141	286267	10.45 ug/ml	99
36) Hexachlorocyclopentadiene	10.22	237	12397	6.62 ug/ml#	84
37) 2,4,6-Trichlorophenol	10.51	196	170938	17.70 ug/ml	99
38) 2,4,5-Trichlorophenol	10.71	196	163777	16.48 ug/ml	98
40) 2-Chloronaphthalene	10.82	162	261273	9.16 ug/ml	99
41) 2-Nitroaniline	11.26	138	116817	8.57 ug/ml	100
42) 2,6-Dinitrotoluene	11.95	165	79164	9.26 ug/ml	100
43) Dimethyl Phthalate	11.79	163	318765	9.45 ug/ml	100
44) Acenaphthylene	11.83	152	428472	9.03 ug/ml	99
45) 3-Nitroaniline	12.28	138	88989	8.28 ug/ml	94
46) Acenaphthene	12.28	154	241086	9.12 ug/ml	97
47) 2,4-dinitrophenol	12.55	184	64829	16.39 ug/ml	95
48) Dibenzofuran	12.67	168	384814	9.61 ug/ml	99
49) 4-Nitrophenol	13.10	109	67805	17.87 ug/ml	95
50) 2,4-Dinitrotoluene	12.89	165	104458	9.57 ug/ml	97
51) 2,3,4,6-Tetrachlorophenol	13.18	232	110991	14.86 ug/ml	98
52) Fluorene	13.50	166	299505	9.74 ug/ml	99
53) Diethyl Phthalate	13.52	149	319865	8.97 ug/ml	98
54) 4-Chlorophenyl phenyl eth	13.56	204	135854	8.79 ug/ml#	89
55) 4-Nitroaniline	13.78	138	95229	8.60 ug/ml	99
56) 4,6-Dinitro-2-methylphenol	13.87	198	123169	17.25 ug/ml	100
57) n-nitrosodiphenylamine	13.91	169	199323	9.56 ug/ml	99
58) Azobenzene	13.93	77	412369	8.47 ug/ml	94
59) Biphenyl	10.83	154	340771	9.14 ug/ml	98
62) 4-bromophenyl phenyl ether	14.72	248	83781	8.77 ug/ml	99
63) Hexachlorobenzene	15.03	284	95817	8.82 ug/ml	96
64) Pentachlorophenol	15.56	266	71871	12.58 ug/ml	93
65) Phenanthrene	15.80	178	396446	9.55 ug/ml	99
66) Anthracene	15.90	178	409770	9.77 ug/ml	99
67) Carbazole	16.39	167	390650	9.70 ug/ml	99
68) Di-n-butyl phthalate	17.44	149	545164	9.04 ug/ml	99
69) Fluoranthene	18.67	202	419518	9.36 ug/ml	99
70) Benzidine	19.12	184	226143	11.96 ug/ml	98

(#) = qualifier out of range (m) = manual integration

Quantitation Report (QT Reviewed)

Data File : D:\HPCHEM\1\DATA\052406-D\D47725SI.D Vial: 4
Acq On : 25 May 2006 4:53 am Operator:
Sample : S6850 PT2 10/20 Inst : D Instrum
Misc : Multiplr: 1.00
Integration Params: RTEINT.P
Quant Time: May 26 9:22 2006 Quant Results File: S05246.RES

Quant Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Fri May 26 09:21:45 2006
Response via : Initial Calibration
DataAcq Meth : ABN

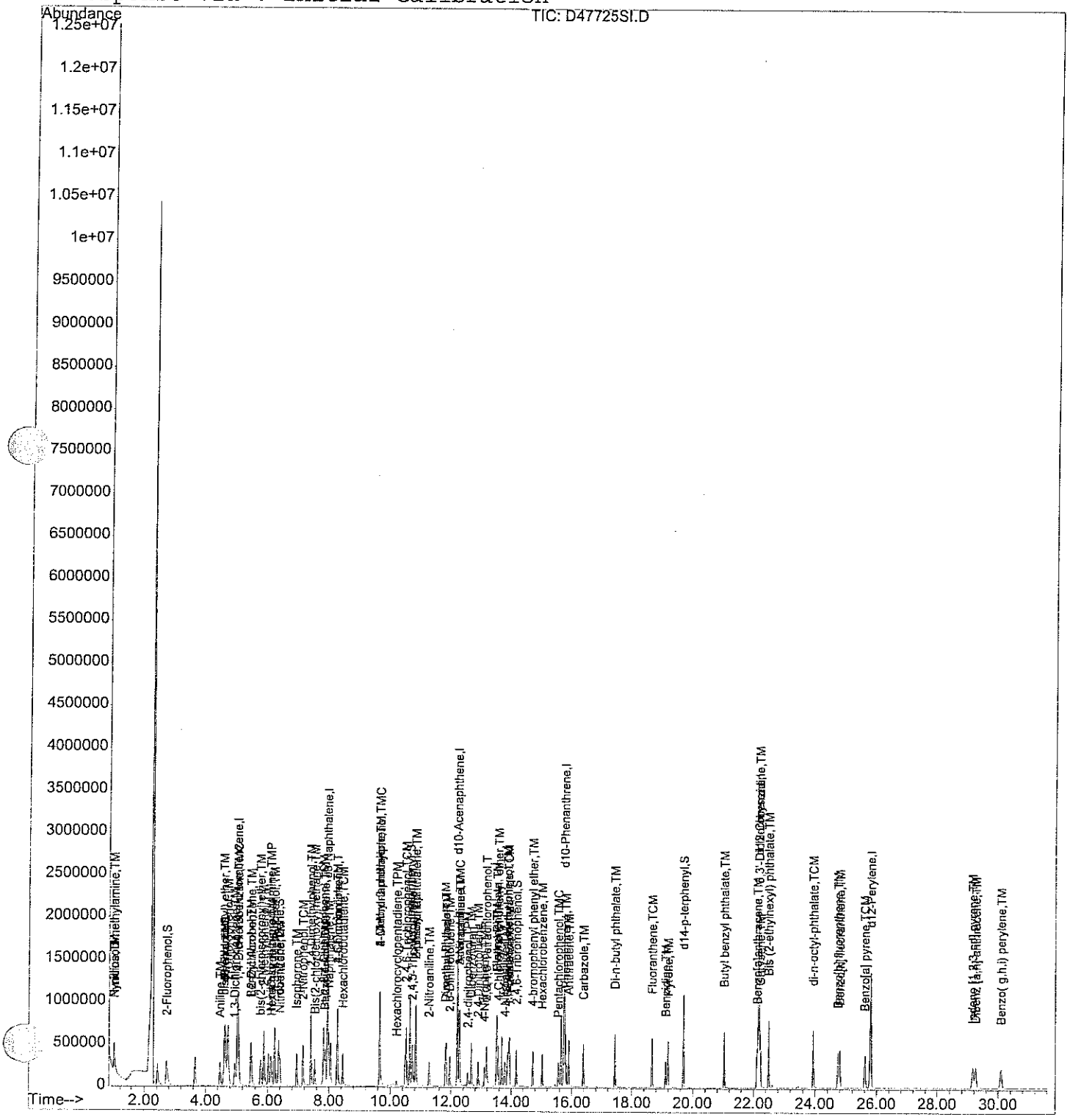
Compound	R.T.	QIon	Response	Conc Unit	Qvalue
72) Pyrene	19.19	202	421297	9.16 ug/ml	99
74) Butyl benzyl phthalate	21.03	149	248667	9.05 ug/ml	99
75) Benzo[a]anthracene	22.11	228	377641	9.20 ug/ml	99
76) Chrysene	22.21	228	352595	9.12 ug/ml	99
77) 3,3'-Dichlorobenzidine	22.18	252	152767	10.35 ug/ml	97
78) Bis (2-ethylhexyl) phthala	22.49	149	338677	9.04 ug/ml	99
80) di-n-octyl-phthalate	23.94	149	592468	9.28 ug/ml	99
81) Benzo[b] fluoranthene	24.73	252	364171	9.26 ug/ml	100
82) Benzo[k] fluoranthene	24.80	252	361533	9.61 ug/ml	99
83) Benzo[a] pyrene	25.60	252	347692	9.32 ug/ml#	92
84) Indeno [1,2,3-cd] pyrene	29.16	276	326017	8.63 ug/ml#	65
85) Dibenz [a,h] anthracene	29.25	278	328015	9.05 ug/ml	99
86) Benzo(g,h,i) perylene	30.10	276	331136	8.92 ug/ml#	85

(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : D:\HPCHEM\1\DATA\052406-D\D47725SI.D Vial: 4
Acq On : 25 May 2006 4:53 am Operator:
Sample : S6850 PT2 10/20 Inst : D Instrum
Misc : Multiplr: 1.00
MS Integration Params: RTEINT.P
Quant Time: May 26 9:22 2006 Quant Results File: S05246.RES

Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Fri May 26 09:35:43 2006
Response via : Initial Calibration



Quantitation Report (QT Reviewed)

Data File : D:\HPCHEM\1\DATA\052406-D\D47726SI.D Vial: 5
Acq On : 25 May 2006 5:26 am Operator:
Sample : S6849 PT3 20/40 Inst : D Instrum
Misc : Multiplr: 1.00
Integration Params: RTEINT.P
Quant Time: May 26 9:22 2006 Quant Results File: S05246.RES

Quant Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Fri May 26 09:21:45 2006
Response via : Initial Calibration
DataAcq Meth : ABN

5.26.06 5.26.06 SL

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) d4-1,4-dichlorobenzene	5.04	152	487489	40.00	ug/ml	0.00
19) d8-Naphthalene	7.95	136	1805489	40.00	ug/ml	0.00
35) d10-Acenaphthene	12.21	164	924277	40.00	ug/ml	0.00
61) d10-Phenanthrene	15.75	188	1566341	40.00	ug/ml	0.00
71) d12-Chrysene	22.17	240	1389789	40.00	ug/ml	0.00
79) d12-Perylene	25.81	264	1317029	40.00	ug/ml	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	2.73	112	763990	43.83	ug/ml	0.00
Spiked Amount 100.000	Range 28 - 110		Recovery = 43.83%			
5) d5-Phenol	4.60	99	912738	44.26	ug/ml	0.00
Spiked Amount 100.000	Range 22 - 110		Recovery = 44.26%			
20) d5-nitrobenzene	6.37	82	827532	40.41	ug/ml	0.00
Spiked Amount 50.000	Range 45 - 107		Recovery = 80.82%			
39) 2-Fluorobiphenyl	10.64	172	1239481	40.55	ug/ml	0.00
Spiked Amount 50.000	Range 44 - 102		Recovery = 81.10%			
60) 2,4,6-Tribromophenol	14.18	330	198286	33.29	ug/ml	0.00
Spiked Amount 100.000	Range 63 - 127		Recovery = 33.29%#			
73) d14-p-terphenyl	19.71	244	1180170	39.43	ug/ml	0.00
Spiked Amount 50.000	Range 33 - 132		Recovery = 78.86%			
Target Compounds						Qvalue
2) Pyridine	1.01	79	388123	14.65	ug/ml#	100
3) N-nitrosodimethylamine	1.00	42	308885	20.17	ug/ml	100
6) Aniline	4.47	93	610787	21.30	ug/ml	100
7) Phenol	4.62	94	976249	43.51	ug/ml	100
8) 2-Chlorophenol	4.73	128	750740	40.49	ug/ml	100
9) bis(2-chloroethyl) ether	4.60	93	395497	20.99	ug/ml	100
10) 1,3-Dichlorobenzene	4.95	146	363791	18.72	ug/ml	100
11) 1,4-Dichlorobenzene	5.07	146	385174	19.51	ug/ml	100
12) 1,2-Dichlorobenzene	5.48	146	358701	19.32	ug/ml	100
13) Benzyl Alcohol	5.47	108	271879	24.84	ug/ml	100
14) 2-Methylphenol	5.88	108	709942	46.01	ug/ml	100
15) bis(2-chloroisopropyl)ethe	5.78	45	612723	16.76	ug/ml	100
16) 3+4-Methylphenol	6.23	108	744744	46.95	ug/ml	100
17) Hexachloroethane	6.13	117	176150	21.58	ug/ml	100
18) N-nitroso-di-n-propylamine	6.11	70	234434	15.73	ug/ml	100
21) Nitrobenzene	6.41	123	210634	21.70	ug/ml	100
22) Isophorone	6.95	82	680195	19.19	ug/ml	100
23) 2-Nitrophenol	7.15	139	411282	38.23	ug/ml	100
24) 2,4-Dimethylphenol	7.41	107	762506	47.99	ug/ml	100
25) Bis(2-chloroethoxy)methane	7.54	93	452563	21.42	ug/ml	100

(#) = qualifier out of range (m) = manual integration

Quantitation Report (QT Reviewed)

Data File : D:\HPCHEM\1\DATA\052406-D\D47726SI.D Vial: 5
Acq On : 25 May 2006 5:26 am Operator:
Sample : S6849 PT3 20/40 Inst : D Instrum
Misc : Multiplr: 1.00
Integration Params: RTEINT.P
Quant Time: May 26 9:22 2006 Quant Results File: S05246.RES

Quant Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Fri May 26 09:21:45 2006
Response via : Initial Calibration
DataAcq Meth : ABN

Compound	R.T.	QIon	Response	Conc Unit	Qvalue
26) 2,4-Dichlorophenol	7.81	162	523537	41.94 ug/ml	100
27) Benzoic Acid	7.93	105	423433	39.42 ug/ml	100
28) 1,2,4-Trichlorobenzene	7.88	180	273686	20.35 ug/ml	100
29) Naphthalene	8.00	128	1000847	20.65 ug/ml	100
30) 2,6-Dichlorophenol	8.29	162	524701	41.83 ug/ml	100
31) 4-Chloroaniline	8.27	127	441817	20.52 ug/ml	100
32) Hexachlorobutadiene	8.46	225	152227	19.39 ug/ml	100
33) 4-Chloro-3-methylphenol	9.65	107	638364	48.63 ug/ml	100
34) 2-Methylnaphthalene	9.66	141	584300	21.28 ug/ml	100
36) Hexachlorocyclopentadiene	10.21	237	49887	14.60 ug/ml	100
37) 2,4,6-Trichlorophenol	10.52	196	353891	36.89 ug/ml	100
38) 2,4,5-Trichlorophenol	10.70	196	355510	36.00 ug/ml	100
40) 2-Chloronaphthalene	10.82	162	542525	19.15 ug/ml	100
41) 2-Nitroaniline	11.26	138	245365	18.11 ug/ml	100
42) 2,6-Dinitrotoluene	11.96	165	164076	19.32 ug/ml	100
43) Dimethyl Phthalate	11.80	163	648737	19.35 ug/ml	100
44) Acenaphthylene	11.83	152	880712	18.67 ug/ml	100
45) 3-Nitroaniline	12.29	138	190254	17.81 ug/ml	100
46) Acenaphthene	12.29	154	487707	18.57 ug/ml	100
47) 2,4-dinitrophenol	12.56	184	173963	35.49 ug/ml	100
48) Dibenzofuran	12.68	168	763596	19.19 ug/ml	100
49) 4-Nitrophenol	13.09	109	143118	37.97 ug/ml	100
50) 2,4-Dinitrotoluene	12.90	165	218127	20.11 ug/ml	100
51) 2,3,4,6-Tetrachlorophenol	13.18	232	238985	32.20 ug/ml	100
52) Fluorene	13.51	166	601438	19.69 ug/ml	100
53) Diethyl Phthalate	13.53	149	656555	18.54 ug/ml	100
54) 4-Chlorophenyl phenyl eth	13.56	204	283244	18.44 ug/ml	100
55) 4-Nitroaniline	13.80	138	199561	18.14 ug/ml	100
56) 4,6-Dinitro-2-methylphenol	13.88	198	274699	35.30 ug/ml	100
57) n-nitrosodiphenylamine	13.91	169	405617	19.57 ug/ml	100
58) Azobenzene	13.94	77	895698	18.52 ug/ml	100
59) Biphenyl	10.84	154	686869	18.54 ug/ml	100
62) 4-bromophenyl phenyl ether	14.72	248	167601	17.68 ug/ml	100
63) Hexachlorobenzene	15.03	284	203588	18.88 ug/ml	100
64) Pentachlorophenol	15.56	266	174419	30.75 ug/ml	100
65) Phenanthrene	15.80	178	783055	19.01 ug/ml	100
66) Anthracene	15.91	178	826397	19.86 ug/ml	100
67) Carbazole	16.40	167	780723	19.52 ug/ml	100
68) Di-n-butyl phthalate	17.45	149	1128654	18.84 ug/ml	100
69) Fluoranthene	18.68	202	832689	18.71 ug/ml	100
70) Benzidine	19.13	184	423906	22.58 ug/ml	100

(#) = qualifier out of range (m) = manual integration

Quantitation Report (QT Reviewed)

Data File : D:\HPCHEM\1\DATA\052406-D\D47726SI.D Vial: 5
Acq On : 25 May 2006 5:26 am Operator:
Sample : S6849 PT3 20/40 Inst : D Instrum
Misc : Multiplr: 1.00
Integration Params: RTEINT.P
Quant Time: May 26 9:22 2006 Quant Results File: S05246.RES

Quant Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Fri May 26 09:21:45 2006
Response via : Initial Calibration
DataAcq Meth : ABN

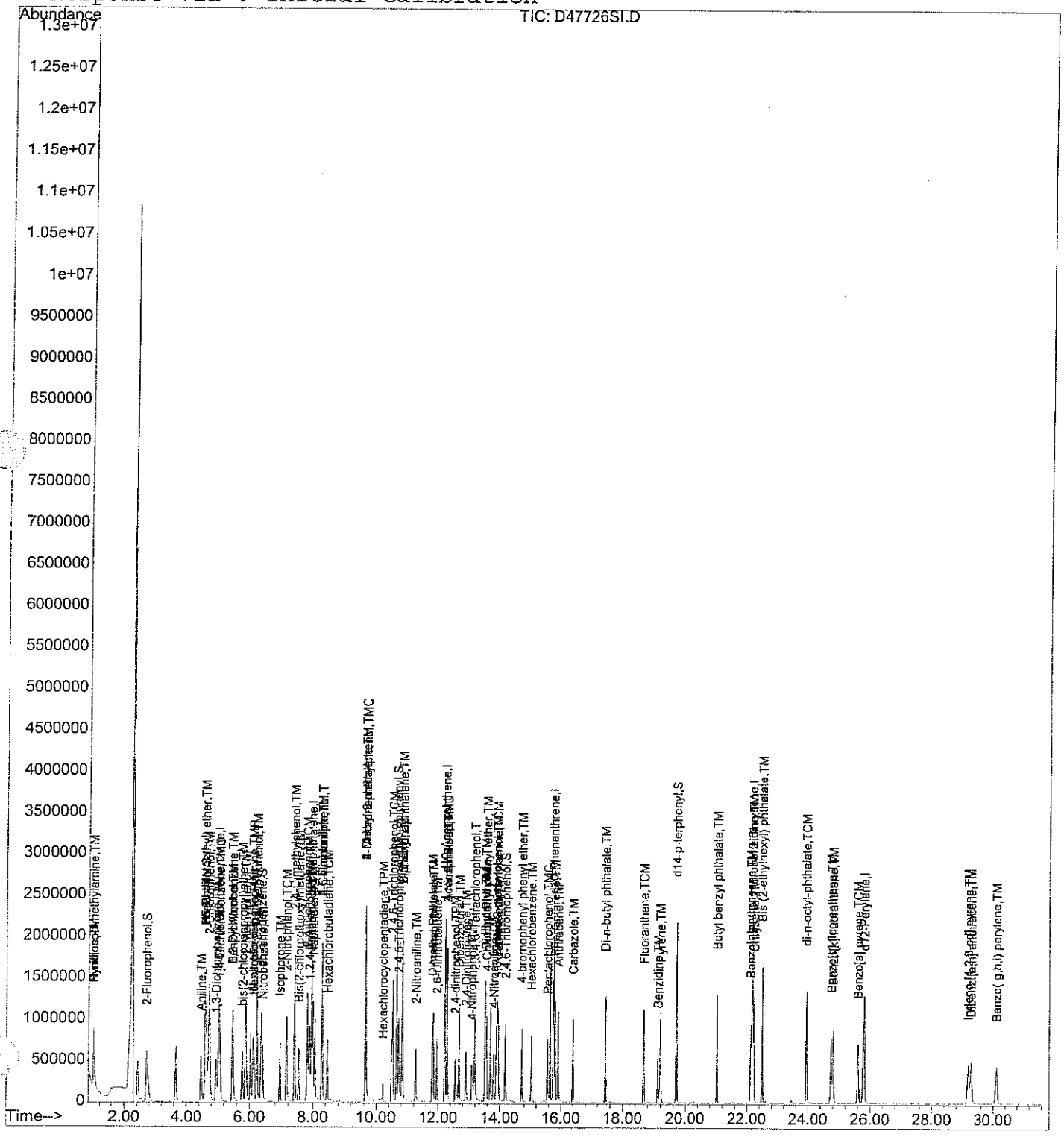
Compound	R.T.	QIon	Response	Conc Unit	Qvalue
72) Pyrene	19.20	202	844614	18.57 ug/ml	100
74) Butyl benzyl phthalate	21.03	149	509016	18.73 ug/ml	100
75) Benzo[a]anthracene	22.12	228	750530	18.48 ug/ml	100
76) Chrysene	22.23	228	713153	18.64 ug/ml	100
77) 3,3'-Dichlorobenzidine	22.19	252	318104	21.79 ug/ml	100
78) Bis (2-ethylhexyl) phthala	22.49	149	690691	18.62 ug/ml	100
80) di-n-octyl-phthalate	23.94	149	1188339	19.02 ug/ml	100
81) Benzo[b] fluoranthene	24.75	252	732023	19.02 ug/ml	99
82) Benzo[k] fluoranthene	24.81	252	716007	19.45 ug/ml	100
83) Benzo[a] pyrene	25.61	252	697882	19.13 ug/ml#	91
84) Indeno [1,2,3-cd] pyrene	29.18	276	672412	18.19 ug/ml#	67
85) Dibenz [a,h] anthracene	29.27	278	660347	18.62 ug/ml	100
86) Benzo(g,h,i) perylene	30.13	276	664919	18.30 ug/ml#	85

(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : D:\HPCHEM\1\DATA\052406-D\D47726SI.D Vial: 5
Acq On : 25 May 2006 5:26 am Operator:
Sample : S6849 PT3 20/40 Inst : D Instrum
Misc : Multiplr: 1.00
MS Integration Params: RTEINT.P
Quant Time: May 26 9:22 2006 Quant Results File: S05246.RES

Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Fri May 26 09:35:43 2006
Response via : Initial Calibration



Quantitation Report (QT Reviewed)

Data File : D:\HPCHEM\1\DATA\052406-D\D47727SI.D Vial: 6
Acq On : 25 May 2006 5:59 am Operator:
Sample : S6848 PT4A 40/80 Inst : D Instrum
Misc : Multiplr: 1.00
Integration Params: RTEINT.P
Quant Time: May 26 9:22 2006 Quant Results File: S05246.RES

Quant Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Fri May 26 09:21:45 2006
Response via : Initial Calibration
DataAcq Meth : ABN

5.26.06 5.26.06 SL

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) d4-1,4-dichlorobenzene	5.04	152	494381	40.00	ug/ml	0.00
19) d8-Naphthalene	7.96	136	1825431	40.00	ug/ml	0.00
35) d10-Acenaphthene	12.22	164	917039	40.00	ug/ml	0.00
61) d10-Phenanthrene	15.76	188	1549650	40.00	ug/ml	0.01
71) d12-Chrysene	22.18	240	1392897	40.00	ug/ml	0.00
79) d12-Perylene	25.81	264	1308524	40.00	ug/ml	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	2.72	112	1562225	88.37	ug/ml	0.00
Spiked Amount 100.000	Range 28 - 110		Recovery = 88.37%			
5) d5-Phenol	4.60	99	1878392	89.82	ug/ml	0.00
Spiked Amount 100.000	Range 22 - 110		Recovery = 89.82%			
20) d5-nitrobenzene	6.39	82	1650899	79.74	ug/ml	0.02
Spiked Amount 50.000	Range 45 - 107		Recovery = 159.48%#			
39) 2-Fluorobiphenyl	10.66	172	2489378	82.09	ug/ml	0.01
Spiked Amount 50.000	Range 44 - 102		Recovery = 164.18%#			
60) 2,4,6-Tribromophenol	14.19	330	420442	71.15	ug/ml	0.01
Spiked Amount 100.000	Range 63 - 127		Recovery = 71.15%			
73) d14-p-terphenyl	19.72	244	2405900	80.19	ug/ml	0.01
Spiked Amount 50.000	Range 33 - 132		Recovery = 160.38%#			
Target Compounds						Qvalue
2) Pyridine	1.01	79	809142	30.12	ug/ml#	100
3) N-nitrosodimethylamine	1.00	42	602335	38.78	ug/ml	97
6) Aniline	4.48	93	1212601	41.70	ug/ml	100
7) Phenol	4.62	94	1996488	87.74	ug/ml	94
8) 2-Chlorophenol	4.74	128	1533049	81.53	ug/ml	99
9) bis(2-chloroethyl) ether	4.62	93	839884	43.95	ug/ml	97
10) 1,3-Dichlorobenzene	4.96	146	743286	37.71	ug/ml	99
11) 1,4-Dichlorobenzene	5.07	146	773168	38.62	ug/ml	98
12) 1,2-Dichlorobenzene	5.47	146	718622	38.16	ug/ml	99
13) Benzyl Alcohol	5.47	108	548891	49.45	ug/ml	99
14) 2-Methylphenol	5.87	108	1432206	91.53	ug/ml	98
15) bis(2-chloroisopropyl)ethe	5.78	45	1212918	32.71	ug/ml	99
16) 3+4-Methylphenol	6.24	108	1503886	93.48	ug/ml	99
17) Hexachloroethane	6.13	117	355380	42.93	ug/ml	99
18) N-nitroso-di-n-propylamine	6.13	70	483004	31.96	ug/ml	99
21) Nitrobenzene	6.42	123	422095	43.01	ug/ml	99
22) Isophorone	6.96	82	1357164	37.87	ug/ml	99
23) 2-Nitrophenol	7.16	139	868880	79.88	ug/ml	96
24) 2,4-Dimethylphenol	7.41	107	1549596	96.46	ug/ml	100
25) Bis(2-chloroethoxy)methane	7.55	93	917875	42.96	ug/ml	100

(#) = qualifier out of range (m) = manual integration

Quantitation Report (QT Reviewed)

Data File : D:\HPCHEM\1\DATA\052406-D\D47727SI.D Vial: 6
Acq On : 25 May 2006 5:59 am Operator:
Sample : S6848 PT4A 40/80 Inst : D Instrum
Misc : Multiplr: 1.00
Integration Params: RTEINT.P
Quant Time: May 26 9:22 2006 Quant Results File: S05246.RES

Quant Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Fri May 26 09:21:45 2006
Response via : Initial Calibration
DataAcq Meth : ABN

Compound	R.T.	QIon	Response	Conc Unit	Qvalue
26) 2,4-Dichlorophenol	7.81	162	1084390	85.91 ug/ml	99
27) Benzoic Acid	8.03	105	1027981	88.56 ug/ml	95
28) 1,2,4-Trichlorobenzene	7.88	180	562431	41.37 ug/ml	100
29) Naphthalene	8.00	128	2026083	41.34 ug/ml	99
30) 2,6-Dichlorophenol	8.30	162	1080368	85.19 ug/ml	100
31) 4-Chloroaniline	8.29	127	911317	41.86 ug/ml	100
32) Hexachlorobutadiene	8.45	225	307825	38.78 ug/ml	99
33) 4-Chloro-3-methylphenol	9.65	107	1303381	98.20 ug/ml	100
34) 2-Methylnaphthalene	9.66	141	1171919	42.22 ug/ml	97
36) Hexachlorocyclopentadiene	10.22	237	148968	33.80 ug/ml	95
37) 2,4,6-Trichlorophenol	10.52	196	726222	76.30 ug/ml	99
38) 2,4,5-Trichlorophenol	10.70	196	756987	77.25 ug/ml	98
40) 2-Chloronaphthalene	10.84	162	1109467	39.47 ug/ml	99
41) 2-Nitroaniline	11.27	138	497340	37.00 ug/ml	99
42) 2,6-Dinitrotoluene	11.97	165	338863	40.21 ug/ml	97
43) Dimethyl Phthalate	11.81	163	1321791	39.74 ug/ml	100
44) Acenaphthylene	11.84	152	1758724	37.58 ug/ml#	69
45) 3-Nitroaniline	12.30	138	397127	37.47 ug/ml	99
46) Acenaphthene	12.29	154	1015574	38.97 ug/ml	98
47) 2,4-dinitrophenol	12.57	184	413367	76.59 ug/ml	99
48) Dibenzofuran	12.69	168	1550125	39.26 ug/ml	100
49) 4-Nitrophenol	13.08	109	304896	81.53 ug/ml	97
50) 2,4-Dinitrotoluene	12.91	165	444751	41.32 ug/ml	98
51) 2,3,4,6-Tetrachlorophenol	13.19	232	518566	70.41 ug/ml	99
52) Fluorene	13.52	166	1228736	40.55 ug/ml	98
53) Diethyl Phthalate	13.55	149	1318544	37.52 ug/ml	100
54) 4-Chlorophenyl phenyl eth	13.57	204	569888	37.39 ug/ml	99
55) 4-Nitroaniline	13.83	138	399331	36.58 ug/ml	99
56) 4,6-Dinitro-2-methylphenol	13.91	198	587653	76.05 ug/ml	98
57) n-nitrosodiphenylamine	13.94	169	818111	39.79 ug/ml	97
58) Azobenzene	13.96	77	1814344	37.82 ug/ml#	82
59) Biphenyl	10.85	154	1416627	38.55 ug/ml	99
62) 4-bromophenyl phenyl ether	14.73	248	345030	36.78 ug/ml	99
63) Hexachlorobenzene	15.04	284	400870	37.57 ug/ml	98
64) Pentachlorophenol	15.57	266	435814	77.66 ug/ml	96
65) Phenanthrene	15.82	178	1585086	38.89 ug/ml	100
66) Anthracene	15.92	178	1590489	38.63 ug/ml	98
67) Carbazole	16.41	167	1576369	39.83 ug/ml	99
68) Di-n-butyl phthalate	17.46	149	2226816	37.58 ug/ml	99
69) Fluoranthene	18.69	202	1710451	38.85 ug/ml	100
70) Benzidine	19.13	184	749079	40.33 ug/ml	99

(#) = qualifier out of range (m) = manual integration

Quantitation Report (QT Reviewed)

Data File : D:\HPCHEM\1\DATA\052406-D\D47727SI.D Vial: 6
Acq On : 25 May 2006 5:59 am Operator:
Sample : S6848 PT4A 40/80 Inst : D Instrum
Misc : Multiplr: 1.00
Integration Params: RTEINT.P
Quant Time: May 26 9:22 2006 Quant Results File: S05246.RES

Quant Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Fri May 26 09:21:45 2006
Response via : Initial Calibration
DataAcq Meth : ABN

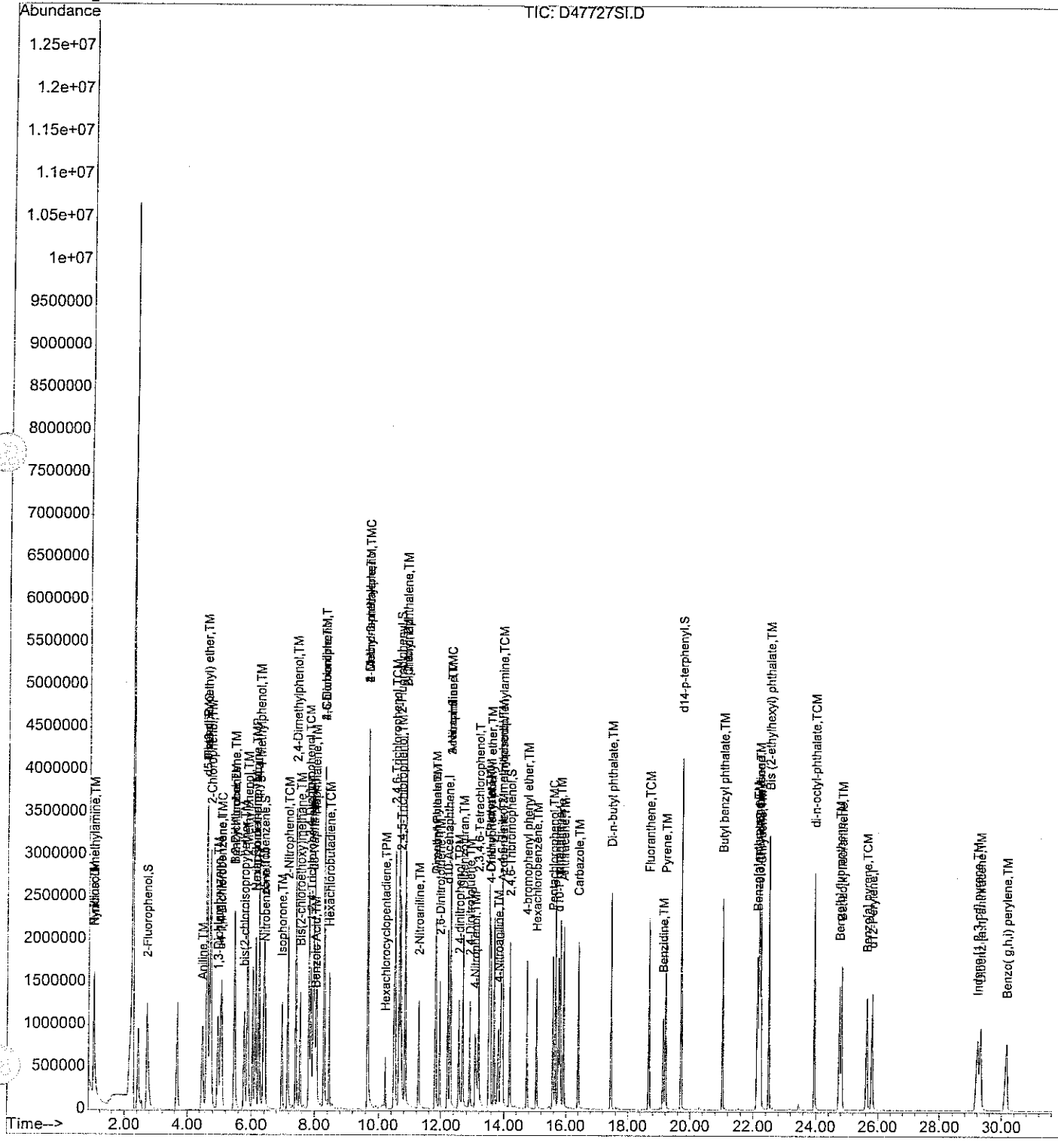
Compound	R.T.	QIon	Response	Conc Unit	Qvalue
72) Pyrene	19.21	202	1713652	37.59 ug/ml	99
74) Butyl benzyl phthalate	21.04	149	1029438	37.80 ug/ml	99
75) Benzo[a]anthracene	22.13	228	1503940	36.94 ug/ml	99
76) Chrysene	22.24	228	1431060	37.31 ug/ml	99
77) 3,3'-Dichlorobenzidine	22.20	252	656817	44.88 ug/ml	99
78) Bis (2-ethylhexyl) phthala	22.50	149	1415781	38.09 ug/ml	99
80) di-n-octyl-phthalate	23.95	149	2421597	39.02 ug/ml	100
81) Benzo[b] fluoranthene	24.76	252	1517895	39.70 ug/ml	99
82) Benzo[k] fluoranthene	24.83	252	1422465	38.90 ug/ml	100
83) Benzo[a] pyrene	25.64	252	1411877	38.95 ug/ml#	91
84) Indeno [1,2,3-cd] pyrene	29.21	276	1424663	38.79 ug/ml#	67
85) Dibenz [a,h] anthracene	29.31	278	1371930	38.93 ug/ml	100
86) Benzo(g,h,i) perylene	30.17	276	1363821	37.79 ug/ml#	85

(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : D:\HPCHEM\1\DATA\052406-D\D47727SI.D Vial: 6
Acq On : 25 May 2006 5:59 am Operator:
Sample : S6848 PT4A 40/80 Inst : D Instrum
Misc : Multiplr: 1.00
MS Integration Params: RTEINT.P
Quant Time: May 26 9:22 2006 Quant Results File: S05246.RES

Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Fri May 26 09:35:43 2006
Response via : Initial Calibration



Quantitation Report (QT Reviewed)

Data File : D:\HPCHEM\1\DATA\052406-D\D47728SI.D Vial: 7
Acq On : 25 May 2006 6:32 am Operator:
Sample : S6847 PT4 50/100 Inst : D Instrum
Misc : Multiplr: 1.00
Integration Params: RTEINT.P
Quant Time: May 26 9:34 2006 Quant Results File: S05246.RES

Quant Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Fri May 26 09:21:45 2006
Response via : Initial Calibration
DataAcq Meth : ABN

5.26.06 5.26.06 SL

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) d4-1,4-dichlorobenzene	5.04	152	486916	40.00	ug/ml	0.00
19) d8-Naphthalene	7.96	136	1808660	40.00	ug/ml	0.00
35) d10-Acenaphthene	12.22	164	934498	40.00	ug/ml	0.00
61) d10-Phenanthrene	15.76	188	1556870	40.00	ug/ml	0.00
71) d12-Chrysene	22.17	240	1376003	40.00	ug/ml	0.00
79) d12-Perylene	25.82	264	1311672	40.00	ug/ml	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	2.72	112	1919998	110.27	ug/ml	0.00
Spiked Amount 100.000	Range 28 - 110		Recovery = 110.27%#			
5) d5-Phenol	4.59	99	2316822	112.49	ug/ml	-0.01
Spiked Amount 100.000	Range 22 - 110		Recovery = 112.49%#			
20) d5-nitrobenzene	6.39	82	2054185	100.14	ug/ml	0.02
Spiked Amount 50.000	Range 45 - 107		Recovery = 200.28%#			
39) 2-Fluorobiphenyl	10.66	172	3098642	100.27	ug/ml	0.02
Spiked Amount 50.000	Range 44 - 102		Recovery = 200.54%#			
60) 2,4,6-Tribromophenol	14.20	330	523687	86.97	ug/ml	0.02
Spiked Amount 100.000	Range 63 - 127		Recovery = 86.97%			
73) d14-p-terphenyl	19.73	244	2996223	101.10	ug/ml	0.02
Spiked Amount 50.000	Range 33 - 132		Recovery = 202.20%#			
Target Compounds						Qvalue
2) Pyridine	1.01	79	989306	37.39	ug/ml#	100
3) N-nitrosodimethylamine	1.00	42	732367	47.87	ug/ml	96
6) Aniline	4.49	93	1505012	52.55	ug/ml	100
7) Phenol	4.63	94	2500608	111.57	ug/ml	94
8) 2-Chlorophenol	4.74	128	1917537	103.54	ug/ml	99
9) bis(2-chloroethyl) ether	4.63	93	1047717	55.67	ug/ml	97
10) 1,3-Dichlorobenzene	4.96	146	923013	47.55	ug/ml	100
11) 1,4-Dichlorobenzene	5.08	146	944979	47.93	ug/ml	98
12) 1,2-Dichlorobenzene	5.48	146	890896	48.03	ug/ml	99
13) Benzyl Alcohol	5.48	108	677077	61.93	ug/ml	97
14) 2-Methylphenol	5.88	108	1777590	115.34	ug/ml	99
15) bis(2-chloroisopropyl)ethe	5.79	45	1520023	41.62	ug/ml	99
16) 3+4-Methylphenol	6.25	108	1877029	118.46	ug/ml	99
17) Hexachloroethane	6.14	117	447294	54.86	ug/ml	99
18) N-nitroso-di-n-propylamine	6.14	70	593018	39.84	ug/ml	98
21) Nitrobenzene	6.43	123	523275	53.81	ug/ml	99
22) Isophorone	6.98	82	1667666	46.96	ug/ml	99
23) 2-Nitrophenol	7.16	139	1075744	99.82	ug/ml	96
24) 2,4-Dimethylphenol	7.42	107	1913980	120.25	ug/ml	99
25) Bis(2-chloroethoxy)methane	7.55	93	1144864	54.08	ug/ml	99

(#) = qualifier out of range (m) = manual integration

Quantitation Report (QT Reviewed)

Data File : D:\HPCHEM\1\DATA\052406-D\D47728SI.D Vial: 7
Acq On : 25 May 2006 6:32 am Operator:
Sample : S6847 PT4 50/100 Inst : D Instrum
Misc : Multiplr: 1.00
Integration Params: RTEINT.P
Quant Time: May 26 9:34 2006 Quant Results File: S05246.RES

Quant Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Fri May 26 09:21:45 2006
Response via : Initial Calibration
DataAcq Meth : ABN

Compound	R.T.	QIon	Response	Conc Unit	Qvalue
26) 2,4-Dichlorophenol	7.81	162	1346488	107.67 ug/ml	99
27) Benzoic Acid	8.07	105	1319032	112.56 ug/ml#	72
28) 1,2,4-Trichlorobenzene	7.89	180	698896	51.88 ug/ml	99
29) Naphthalene	8.01	128	2490853	51.30 ug/ml	99
30) 2,6-Dichlorophenol	8.30	162	1346017	107.13 ug/ml	100
31) 4-Chloroaniline	8.30	127	1147613	53.20 ug/ml	99
32) Hexachlorobutadiene	8.46	225	383623	48.77 ug/ml	98
33) 4-Chloro-3-methylphenol	9.65	107	1656662	125.97 ug/ml	99
34) 2-Methylnaphthalene	9.67	141	1473460	53.57 ug/ml	98
36) Hexachlorocyclopentadiene	10.22	237	201752	42.39 ug/ml	96
37) 2,4,6-Trichlorophenol	10.53	196	900989	92.89 ug/ml	98
38) 2,4,5-Trichlorophenol	10.70	196	924217	92.56 ug/ml	96
40) 2-Chloronaphthalene	10.84	162	1386032	48.39 ug/ml	99
41) 2-Nitroaniline	11.28	138	615728	44.95 ug/ml	99
42) 2,6-Dinitrotoluene	11.97	165	423329	49.29 ug/ml	96
43) Dimethyl Phthalate	11.81	163	1633987	48.21 ug/ml	100
44) Acenaphthylene	11.84	152	2186829	45.86 ug/ml#	70
45) 3-Nitroaniline	12.30	138	498280	46.13 ug/ml	97
46) Acenaphthene	12.30	154	1259628	47.43 ug/ml	98
47) 2,4-dinitrophenol	12.57	184	535894	95.25 ug/ml	99
48) Dibenzofuran	12.69	168	1921281	47.76 ug/ml	100
49) 4-Nitrophenol	13.07	109	396622	104.07 ug/ml	100
50) 2,4-Dinitrotoluene	12.92	165	556053	50.70 ug/ml	99
51) 2,3,4,6-Tetrachlorophenol	13.20	232	650550	86.68 ug/ml	99
52) Fluorene	13.53	166	1520907	49.26 ug/ml	98
53) Diethyl Phthalate	13.55	149	1648371	46.03 ug/ml	99
54) 4-Chlorophenyl phenyl eth	13.58	204	709117	45.65 ug/ml	99
55) 4-Nitroaniline	13.85	138	483285m	43.44 ug/ml	
56) 4,6-Dinitro-2-methylphenol	13.92	198	737876	95.30 ug/ml	98
57) n-nitrosodiphenylamine	13.94	169	1012699	48.33 ug/ml	96
58) Azobenzene	13.96	77	2212925	45.26 ug/ml#	81
59) Biphenyl	10.85	154	1754809	46.86 ug/ml	100
62) 4-bromophenyl phenyl ether	14.73	248	434956	46.15 ug/ml	98
63) Hexachlorobenzene	15.04	284	502796	46.91 ug/ml	98
64) Pentachlorophenol	15.57	266	545761	96.80 ug/ml	96
65) Phenanthrene	15.82	178	1954203	47.72 ug/ml	100
66) Anthracene	15.93	178	1987536	48.04 ug/ml	99
67) Carbazole	16.41	167	1890488	47.55 ug/ml	100
68) Di-n-butyl phthalate	17.46	149	2750634	46.20 ug/ml	100
69) Fluoranthene	18.69	202	2102295	47.53 ug/ml	100
70) Benzidine	19.13	184	903849	48.44 ug/ml	99

(#) = qualifier out of range (m) = manual integration

Quantitation Report (QT Reviewed)

Data File : D:\HPCHEM\1\DATA\052406-D\D47728SI.D Vial: 7
Acq On : 25 May 2006 6:32 am Operator:
Sample : S6847 PT4 50/100 Inst : D Instrum
Misc : Multiplr: 1.00
Integration Params: RTEINT.P
Quant Time: May 26 9:34 2006 Quant Results File: S05246.RES

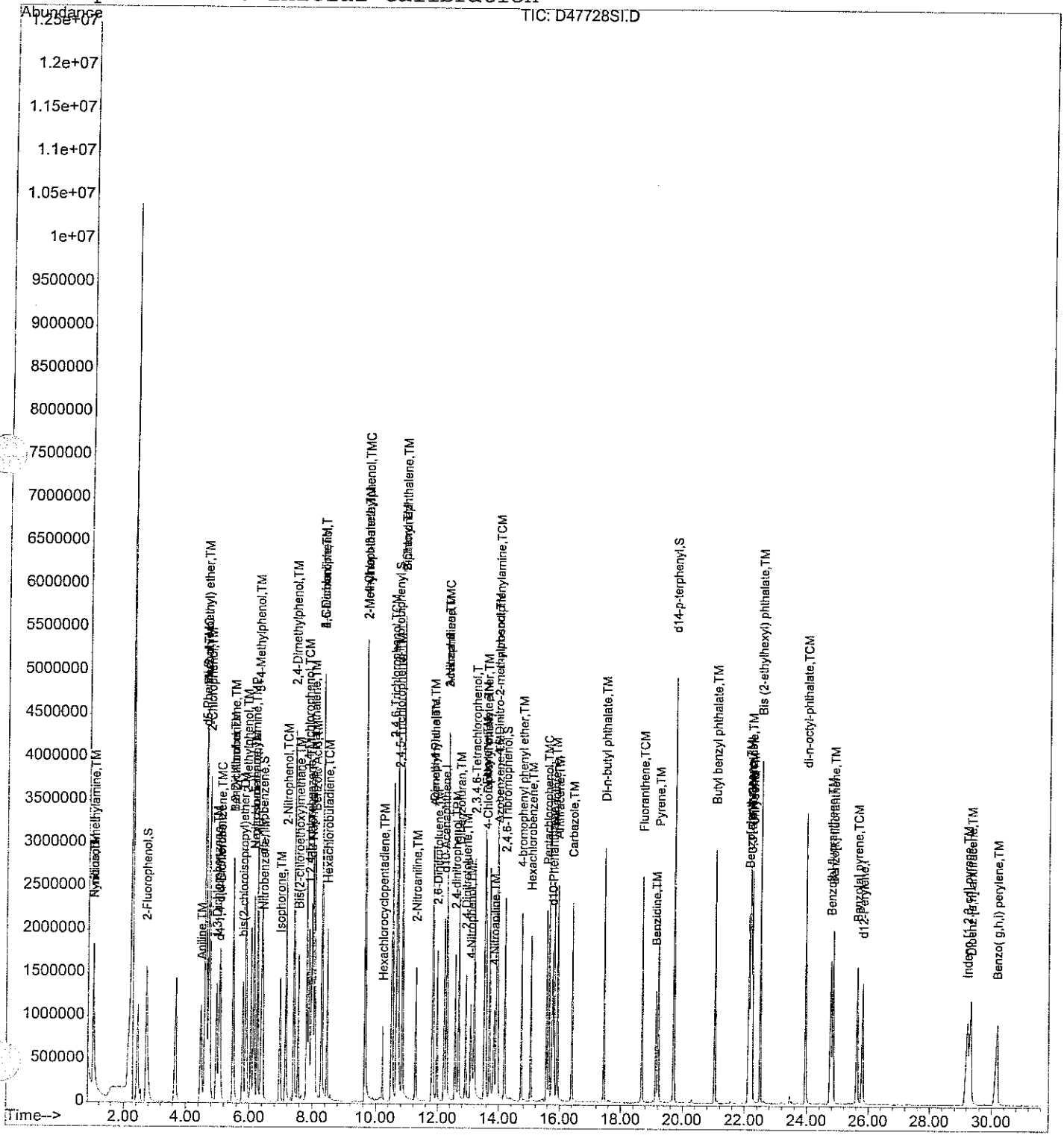
Quant Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Fri May 26 09:21:45 2006
Response via : Initial Calibration
DataAcq Meth : ABN

Compound	R.T.	QIon	Response	Conc Unit	Qvalue
72) Pyrene	19.22	202	2094273	46.50 ug/ml	99
74) Butyl benzyl phthalate	21.04	149	1254300	46.62 ug/ml	99
75) Benzo[a]anthracene	22.13	228	1862989	46.33 ug/ml	99
76) Chrysene	22.24	228	1758491	46.41 ug/ml	98
77) 3,3'-Dichlorobenzidine	22.21	252	813548	56.28 ug/ml	99
78) Bis (2-ethylhexyl) phthala	22.50	149	1753758	47.76 ug/ml	98
80) di-n-octyl-phthalate	23.96	149	3001021	48.24 ug/ml	100
81) Benzo[b] fluoranthene	24.77	252	1845686	48.16 ug/ml	99
82) Benzo[k] fluoranthene	24.84	252	1770217	48.29 ug/ml	100
83) Benzo[a] pyrene	25.65	252	1744913	48.02 ug/ml#	92
84) Indeno [1,2,3-cd] pyrene	29.23	276	1833139	49.80 ug/ml#	67
85) Dibenz [a,h] anthracene	29.33	278	1764541	49.95 ug/ml	98
86) Benzo(g,h,i) perylene	30.19	276	1746121	48.27 ug/ml#	85

(#) = qualifier out of range (m) = manual integration

Data File : D:\HPCHEM\1\DATA\052406-D\D47728SI.D Vial: 7
Acq On : 25 May 2006 6:32 am Operator:
Sample : S6847 PT4 50/100 Inst : D Instrum
Misc : Multiplr: 1.00
MS Integration Params: RTEINT.P
Quant Time: May 26 9:34 2006 Quant Results File: S05246.RES

Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Fri May 26 09:35:43 2006
Response via : Initial Calibration



Quantitation Report (Qedit)

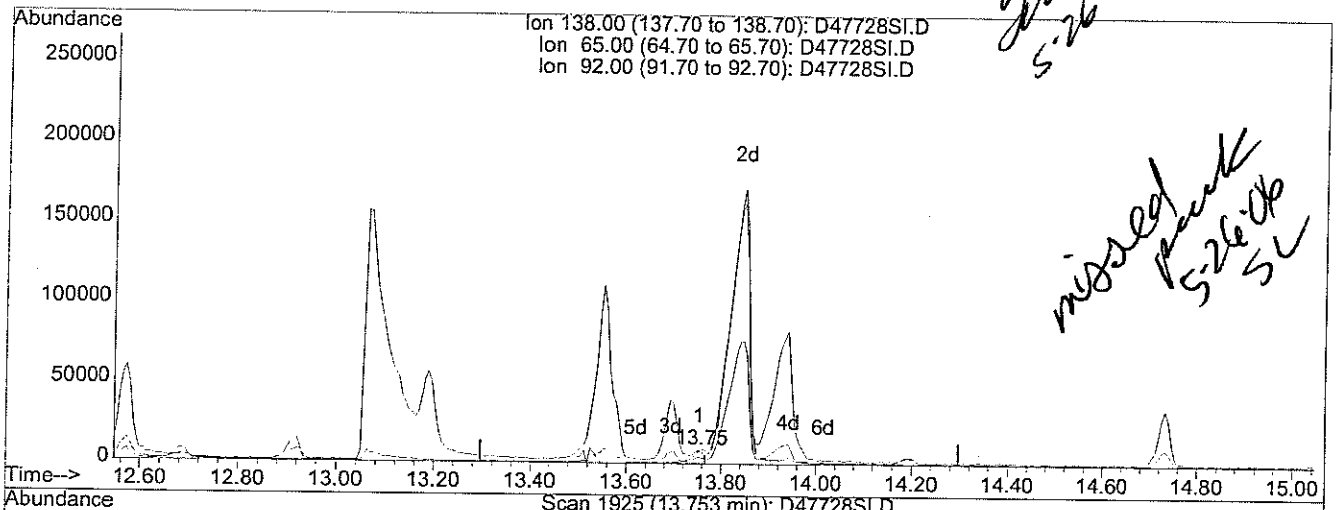
Data File : D:\HPCHEM\1\DATA\052406-D\D47728SI.D
Acq On : 25 May 2006 6:32 am
Sample : S6847 PT4 50/100
Misc :
MS Integration Params: RTEINT.P
Quant Time: May 26 9:22 2006

Vial: 7
Operator:
Inst : D Instrum
Multiplr: 1.00

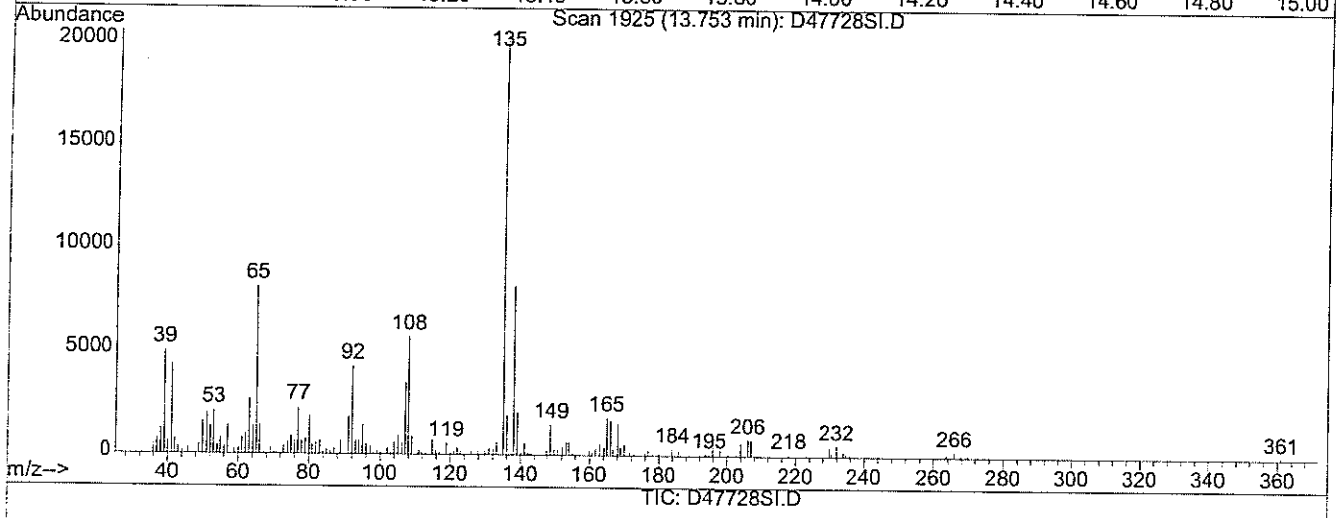
Quant Results File: temp.res

Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Fri May 26 09:30:40 2006
Response via : Multiple Level Calibration

5-26-06



missed peak 5-26-06 SL



(55) 4-Nitroaniline (TM)

13.75min 1.32ug/ml

response 14693

Ion	Exp%	Act%
138.00	100	100
65.00	96.40	0.00#
92.00	47.80	56.65
0.00	0.00	0.00

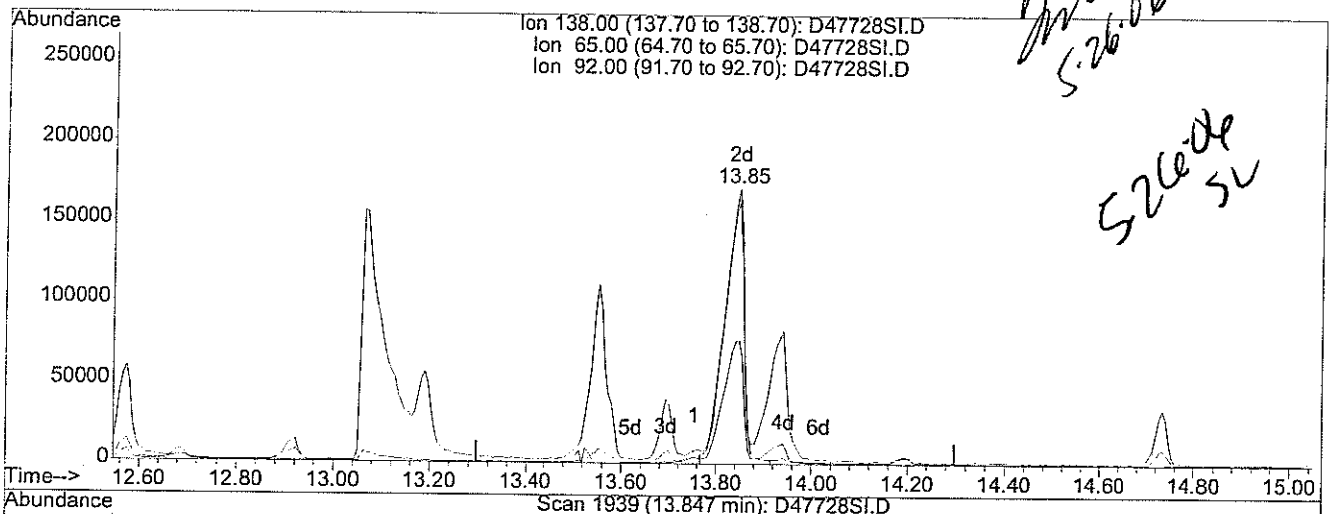
Quantitation Report (Qedit)

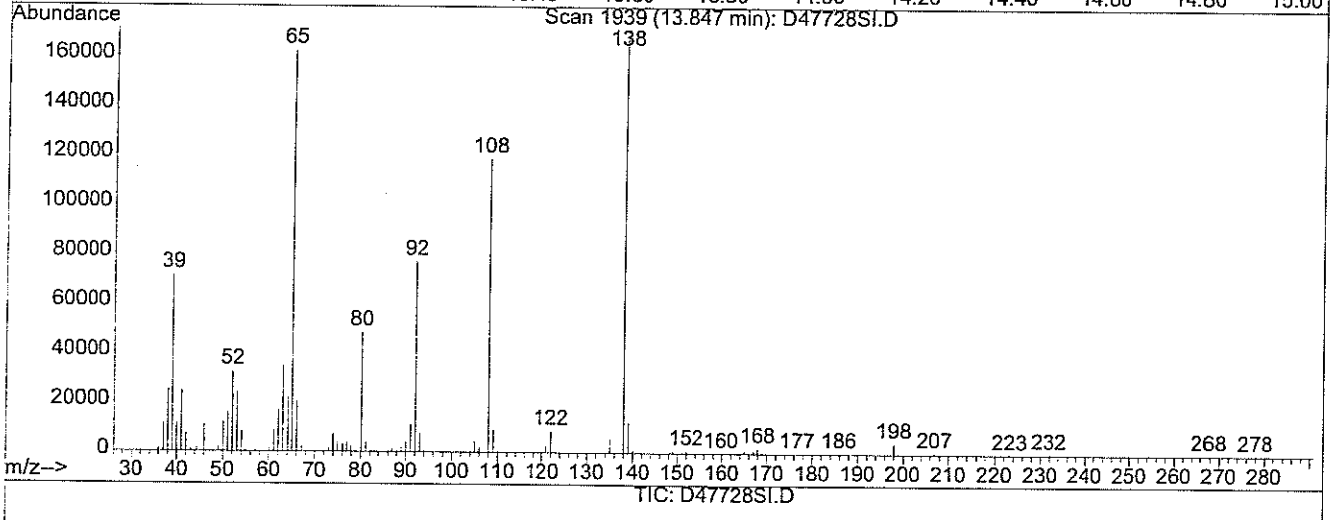
Data File : D:\HPCHEM\1\DATA\052406-D\D47728SI.D
Acq On : 25 May 2006 6:32 am
Sample : S6847 PT4 50/100
Misc :
MS Integration Params: RTEINT.P
Quant Time: May 26 9:34 2006

Vial: 7
Operator:
Inst : D Instrum
Multiplr: 1.00

Quant Results File: temp.res

Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Fri May 26 09:30:40 2006
Response via : Multiple Level Calibration





(55) 4-Nitroaniline (TM)

13.85min 43.44ug/ml m

response 483285

Ion	Exp%	Act%
138.00	100	100
65.00	96.40	0.00#
92.00	47.80	1.72#
0.00	0.00	0.00

Quantitation Report (QT Reviewed)

Data File : D:\HPCHEM\1\DATA\052406-D\D47729SI.D Vial: 8
Acq On : 25 May 2006 7:06 am Operator:
Sample : S6846 PT5 100/200 Inst : D Instrum
Misc : Multiplr: 1.00
Integration Params: RTEINT.P
Quant Time: May 26 9:29 2006 Quant Results File: S05246.RES

Quant Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Fri May 26 09:21:45 2006
Response via : Initial Calibration
DataAcq Meth : ABN

5.26.06 5.26.06 SC

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) d4-1,4-dichlorobenzene	5.05	152	502892	40.00	ug/ml	0.00
19) d8-Naphthalene	7.98	136	1847600	40.00	ug/ml	0.03
35) d10-Acenaphthene	12.23	164	942480	40.00	ug/ml	0.01
61) d10-Phenanthrene	15.77	188	1557195	40.00	ug/ml	0.02
71) d12-Chrysene	22.19	240	1196579	40.00	ug/ml	0.01
79) d12-Perylene	25.83	264	1227072	40.00	ug/ml	0.02
System Monitoring Compounds						
4) 2-Fluorophenol	2.72	112	3887405	216.18	ug/ml	0.00
Spiked Amount 100.000	Range 28 - 110		Recovery = 216.18%#			
5) d5-Phenol	4.62	99	4851590	228.07	ug/ml	0.01
Spiked Amount 100.000	Range 22 - 110		Recovery = 228.07%#			
20) d5-nitrobenzene	6.42	82	4095524	195.45	ug/ml	0.06
Spiked Amount 50.000	Range 45 - 107		Recovery = 390.90%#			
39) 2-Fluorobiphenyl	10.70	172	6802770m	218.27	ug/ml	0.06
Spiked Amount 50.000	Range 44 - 102		Recovery = 436.54%#			
60) 2,4,6-Tribromophenol	14.21	330	1120337	184.48	ug/ml	0.04
Spiked Amount 100.000	Range 63 - 127		Recovery = 184.48%#			
73) d14-p-terphenyl	19.74	244	5702540	221.26	ug/ml	0.04
Spiked Amount 50.000	Range 33 - 132		Recovery = 442.52%#			
Target Compounds						Qvalue
2) Pyridine	1.02	79	1997857	73.10	ug/ml#	100
3) N-nitrosodimethylamine	1.01	42	1477983	93.55	ug/ml	96
6) Aniline	4.50	93	3203623	108.31	ug/ml	96
7) Phenol	4.65	94	5189444	224.19	ug/ml	89
8) 2-Chlorophenol	4.77	128	3965860	207.34	ug/ml	99
9) bis(2-chloroethyl) ether	4.67	93	1905596	98.03	ug/ml	91
10) 1,3-Dichlorobenzene	4.97	146	1851601	92.36	ug/ml	99
11) 1,4-Dichlorobenzene	5.09	146	1924992	94.53	ug/ml	100
12) 1,2-Dichlorobenzene	5.48	146	1805136	94.22	ug/ml	98
13) Benzyl Alcohol	5.49	108	1403114	124.26	ug/ml	99
14) 2-Methylphenol	5.88	108	3638447	228.59	ug/ml	99
15) bis(2-chloroisopropyl)ethe	5.79	45	2952323	78.28	ug/ml	98
16) 3+4-Methylphenol	6.28	108	3889499	237.67	ug/ml	98
17) Hexachloroethane	6.14	117	899502	106.82	ug/ml	99
18) N-nitroso-di-n-propylamine	6.19	70	1202508m	78.21	ug/ml	
21) Nitrobenzene	6.46	123	1070063	107.72	ug/ml	98
22) Isophorone	7.03	82	3325385m	91.67	ug/ml	
23) 2-Nitrophenol	7.18	139	2177864	197.82	ug/ml	95
24) 2,4-Dimethylphenol	7.44	107	3880323	238.65	ug/ml	98
25) Bis(2-chloroethoxy)methane	7.58	93	2247453	103.92	ug/ml	99

(#) = qualifier out of range (m) = manual integration

Quantitation Report (QT Reviewed)

Data File : D:\HPCHEM\1\DATA\052406-D\D47729SI.D Vial: 8
Acq On : 25 May 2006 7:06 am Operator:
Sample : S6846 PT5 100/200 Inst : D Instrum
Misc : Multiplr: 1.00
Integration Params: RTEINT.P
Quant Time: May 26 9:29 2006 Quant Results File: S05246.RES

Quant Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Fri May 26 09:21:45 2006
Response via : Initial Calibration
DataAcq Meth : ABN

Compound	R.T.	QIon	Response	Conc Unit	Qvalue
26) 2,4-Dichlorophenol	7.83	162	2784213	217.94 ug/ml	97
27) Benzoic Acid	8.21	105	2859129m	223.77 ug/ml	
28) 1,2,4-Trichlorobenzene	7.90	180	1425375	103.58 ug/ml	100
29) Naphthalene	8.03	128	4954222	99.88 ug/ml	98
30) 2,6-Dichlorophenol	8.32	162	2929416	228.23 ug/ml	93
31) 4-Chloroaniline	8.32	127	2494801	113.22 ug/ml	96
32) Hexachlorobutadiene	8.47	225	795357	98.99 ug/ml	99
33) 4-Chloro-3-methylphenol	9.66	107	3545027	263.88 ug/ml	98
34) 2-Methylnaphthalene	9.68	141	3222482	114.69 ug/ml	98
36) Hexachlorocyclopentadiene	10.22	237	518208	87.91 ug/ml	97
37) 2,4,6-Trichlorophenol	10.54	196	1900823	194.31 ug/ml	98
38) 2,4,5-Trichlorophenol	10.71	196	2406403	238.96 ug/ml	96
40) 2-Chloronaphthalene	10.87	162	3209997	111.13 ug/ml	99
41) 2-Nitroaniline	11.30	138	1248001	90.33 ug/ml	98
42) 2,6-Dinitrotoluene	12.00	165	840763	97.07 ug/ml	97
43) Dimethyl Phthalate	11.85	163	3367160	98.51 ug/ml	97
44) Acenaphthylene	11.85	152	4410874	91.71 ug/ml#	69
45) 3-Nitroaniline	12.33	138	1012589	92.95 ug/ml	96
46) Acenaphthene	12.32	154	2575442	96.15 ug/ml	98
47) 2,4-dinitrophenol	12.60	184	1191292	194.77 ug/ml	98
48) Dibenzofuran	12.71	168	3844807	94.76 ug/ml	100
49) 4-Nitrophenol	13.06	109	848523	220.76 ug/ml	98
50) 2,4-Dinitrotoluene	12.95	165	1135022	102.61 ug/ml	98
51) 2,3,4,6-Tetrachlorophenol	13.21	232	1425340	188.31 ug/ml	99
52) Fluorene	13.54	166	3063650	98.39 ug/ml	99
53) Diethyl Phthalate	13.58	149	2877020m	79.65 ug/ml	
54) 4-Chlorophenyl phenyl eth	13.59	204	1575950	100.60 ug/ml	95
55) 4-Nitroaniline	13.90	138	822653m	73.32 ug/ml	
56) 4,6-Dinitro-2-methylphenol	13.96	198	1449894m	219.23 ug/ml	
57) n-nitrosodiphenylamine	13.97	169	2220590m	105.08 ug/ml	
58) Azobenzene	13.98	77	4539612	92.07 ug/ml#	81
59) Biphenyl	10.87	154	4080405	108.04 ug/ml	98
62) 4-bromophenyl phenyl ether	14.74	248	869397	92.23 ug/ml	98
63) Hexachlorobenzene	15.06	284	1035804	96.61 ug/ml	99
64) Pentachlorophenol	15.59	266	1176651	208.65 ug/ml	94
65) Phenanthrene	15.84	178	4126381	100.75 ug/ml	100
66) Anthracene	15.95	178	3653236	88.29 ug/ml	98
67) Carbazole	16.43	167	3641294	91.57 ug/ml	97
68) Di-n-butyl phthalate	17.47	149	4511141	75.75 ug/ml	99
69) Fluoranthene	18.70	202	3708547	83.84 ug/ml	96
70) Benzidine	19.14	184	1301158	69.72 ug/ml	99

(#) = qualifier out of range (m) = manual integration

Quantitation Report (QT Reviewed)

Data File : D:\HPCHEM\1\DATA\052406-D\D47729SI.D Vial: 8
Acq On : 25 May 2006 7:06 am Operator:
Sample : S6846 PT5 100/200 Inst : D Instrum
Misc : Multiplr: 1.00
Integration Params: RTEINT.P
Quant Time: May 26 9:29 2006 Quant Results File: S05246.RES

Quant Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Fri May 26 09:21:45 2006
Response via : Initial Calibration
DataAcq Meth : ABN

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
72) Pyrene	19.23	202	3567096	91.08	ug/ml	98
74) Butyl benzyl phthalate	21.06	149	2200737	94.06	ug/ml	99
75) Benzo[a]anthracene	22.15	228	3164703	90.50	ug/ml	98
76) Chrysene	22.27	228	3113194	94.48	ug/ml	99
77) 3,3'-Dichlorobenzidine	22.22	252	1484179	118.06	ug/ml	98
78) Bis (2-ethylhexyl) phthala	22.52	149	3396115	106.36	ug/ml	97
80) di-n-octyl-phthalate	23.98	149	5508414	94.65	ug/ml	100
81) Benzo[b] fluoranthene	24.80	252	3368715	93.97	ug/ml	99
82) Benzo[k] fluoranthene	24.88	252	3194200m	93.14	ug/ml	
83) Benzo[a] pyrene	25.68	252	3179753	93.54	ug/ml#	91
84) Indeno [1,2,3-cd] pyrene	29.28	276	3652406	106.06	ug/ml#	65
85) Dibenz [a,h] anthracene	29.41	278	3604129	109.05	ug/ml	99
86) Benzo(g,h,i) perylene	30.27	276	3406455m	100.65	ug/ml	

(#) = qualifier out of range (m) = manual integration

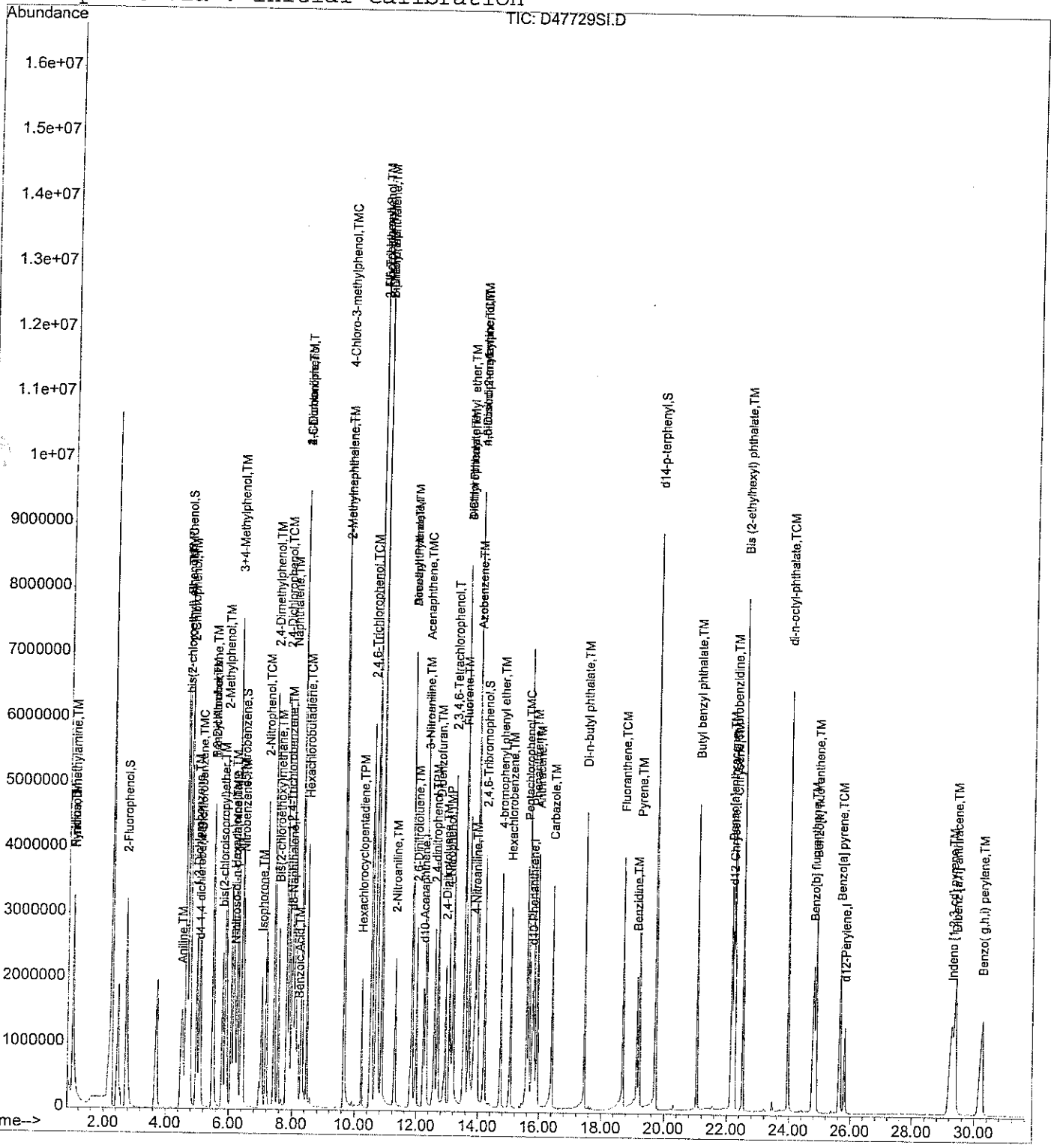
Quantitation Report

Data File : D:\HPCHEM\1\DATA\052406-D\D47729SI.D
Acq On : 25 May 2006 7:06 am
Sample : S6846 PT5 100/200
Misc :
MS Integration Params: RTEINT.P
Quant Time: May 26 9:29 2006

Vial: 8
Operator:
Inst : D Instrum
Multiplr: 1.00

Quant Results File: S05246.RES

Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Fri May 26 09:35:43 2006
Response via : Initial Calibration



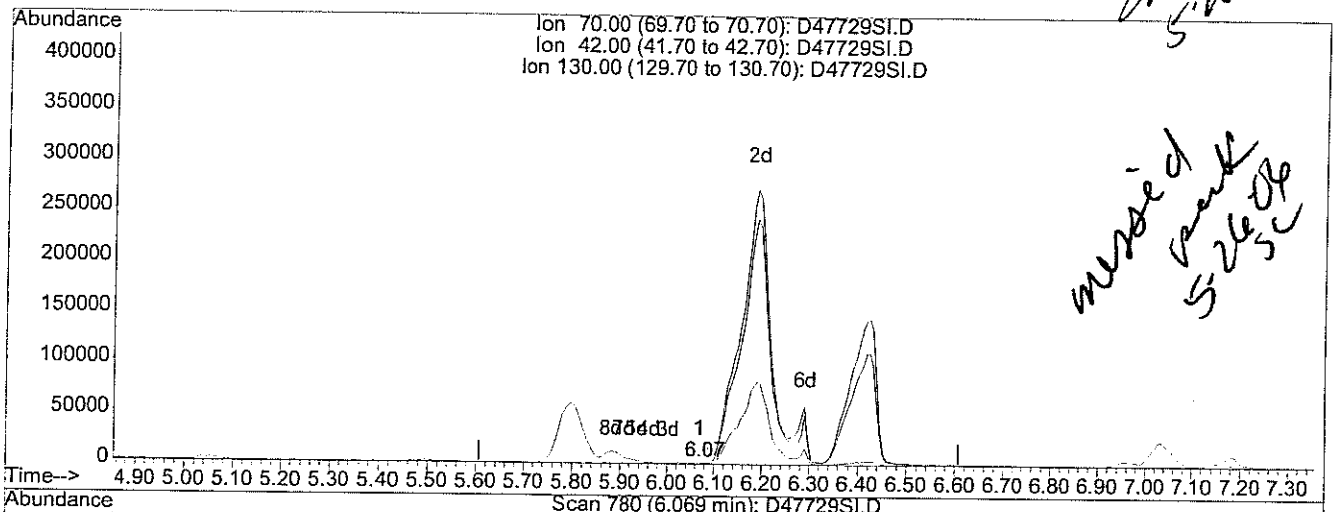
Quantitation Report (Qedit)

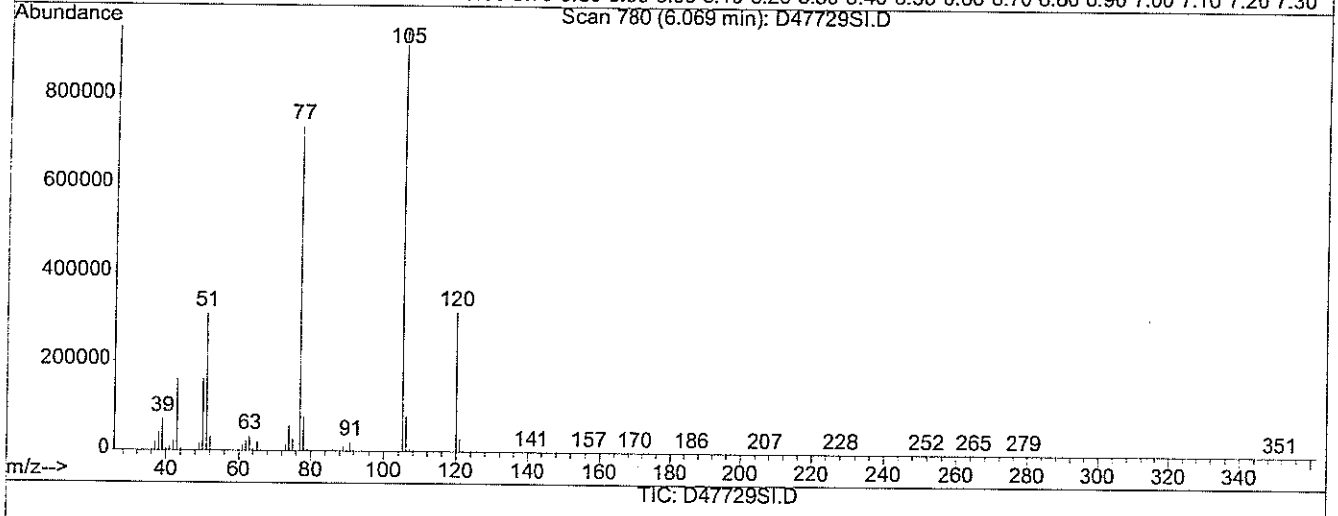
Data File : D:\HPCHEM\1\DATA\052406-D\D47729SI.D
Acq On : 25 May 2006 7:06 am
Sample : S6846 PT5 100/200
Misc :
MS Integration Params: RTEINT.P
Quant Time: May 26 9:22 2006

Vial: 8
Operator:
Inst : D Instrum
Multiplr: 1.00

Quant Results File: temp.res

Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Fri May 26 09:21:45 2006
Response via : Multiple Level Calibration





TIC: D47729SI.D

(18) N-nitroso-di-n-propylamine (TMP)

6.07min 0.25ug/ml

response 3920

Ion	Exp%	Act%
70.00	100	100
42.00	87.70	1556.61#
130.00	31.10	2.12#
0.00	0.00	0.00

Quantitation Report (Qedit)

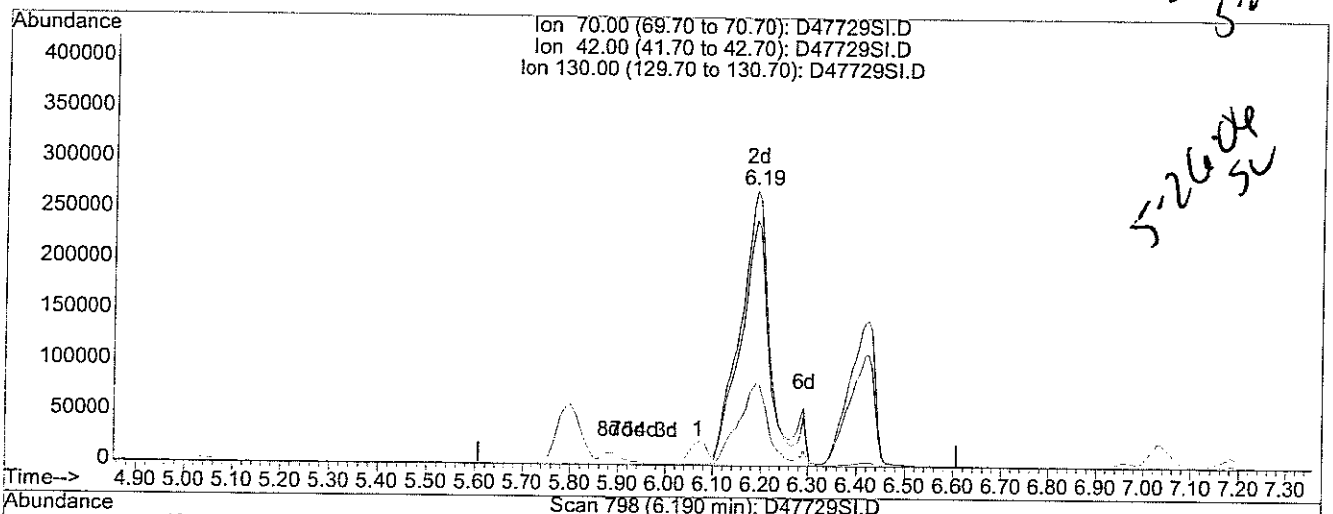
Data File : D:\HPCHEM\1\DATA\052406-D\D47729SI.D
Acq On : 25 May 2006 7:06 am
Sample : S6846 PT5 100/200
Misc :
MS Integration Params: RTEINT.P
Quant Time: May 26 9:27 2006

Vial: 8
Operator:
Inst : D Instrum
Multiplr: 1.00

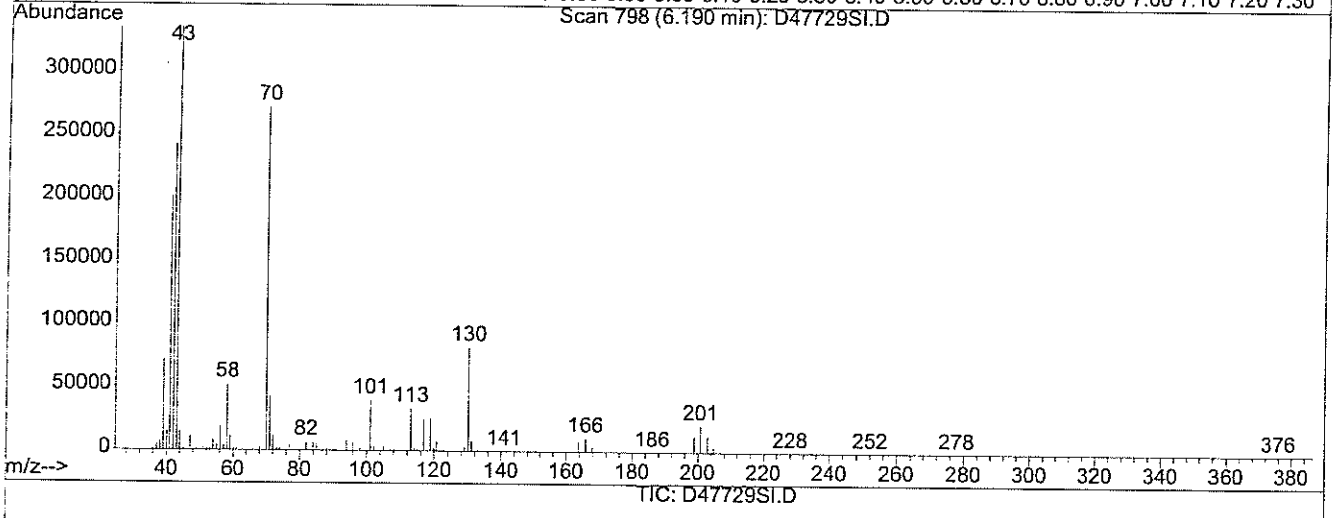
Quant Results File: temp.res

Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Fri May 26 09:21:45 2006
Response via : Multiple Level Calibration

5.26.06



5.26.06 SC



(18) N-nitroso-di-n-propylamine (TMP)

6.19min 78.21ug/ml m

response 1202508

Ion	Exp%	Act%
70.00	100	100
42.00	87.70	5.07#
130.00	31.10	0.01#
0.00	0.00	0.00

Quantitation Report (Qedit)

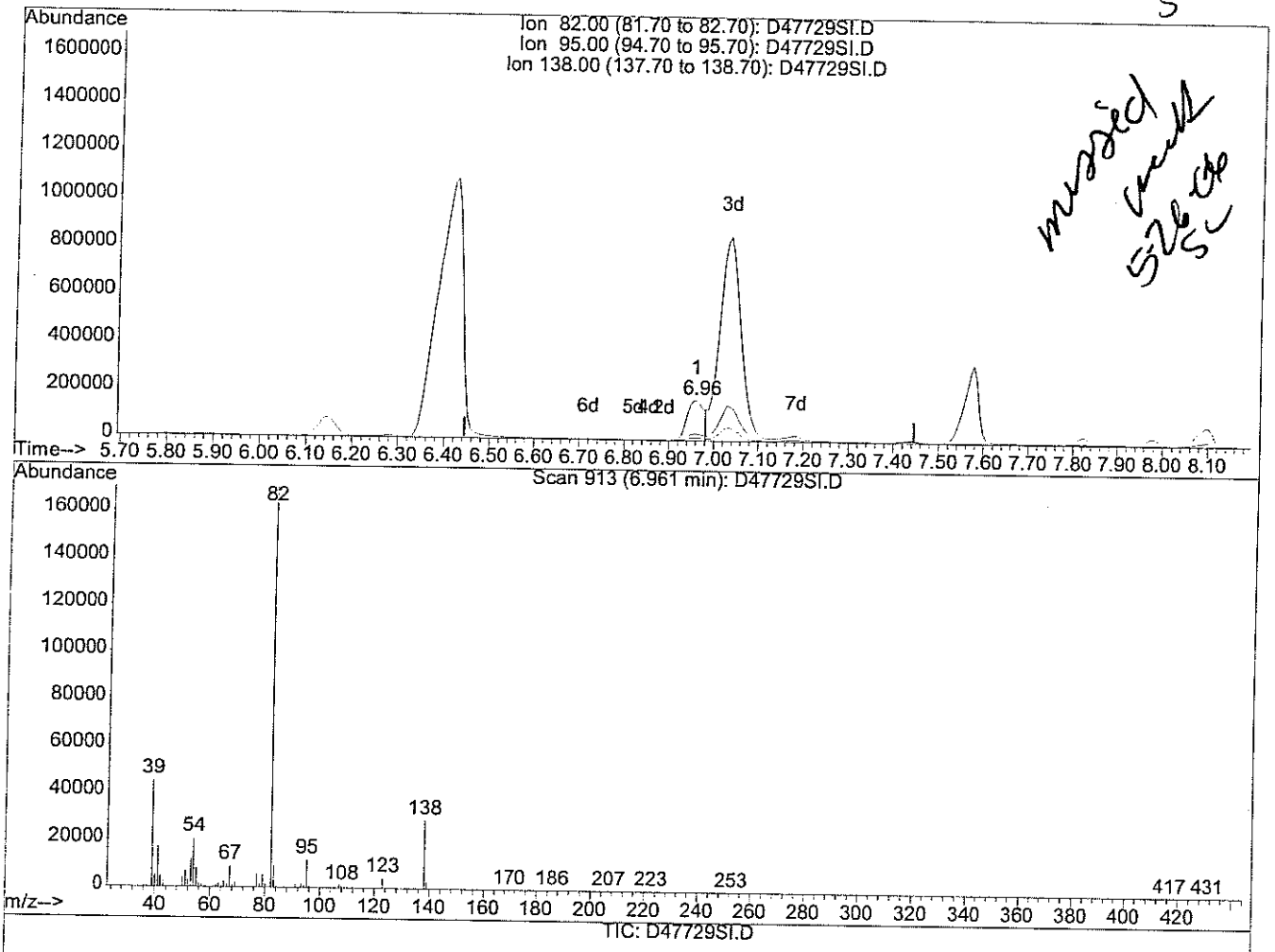
Data File : D:\HPCHEM\1\DATA\052406-D\D47729SI.D
Acq On : 25 May 2006 7:06 am
Sample : S6846 PT5 100/200
Misc :
MS Integration Params: RTEINT.P
Quant Time: May 26 9:27 2006

Vial: 8
Operator:
Inst : D Instrum
Multiplr: 1.00

Quant Results File: temp.res

Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Fri May 26 09:21:45 2006
Response via : Multiple Level Calibration

5.26.06



missed result 5-26-06

TIC: D47729SI.D

(22) Isophorone (TM)

6.96min 12.25ug/ml

response 444446

Ion	Exp%	Act%
82.00	100	100
95.00	6.20	7.88#
138.00	16.50	16.84
0.00	0.00	0.00

Quantitation Report (Qedit)

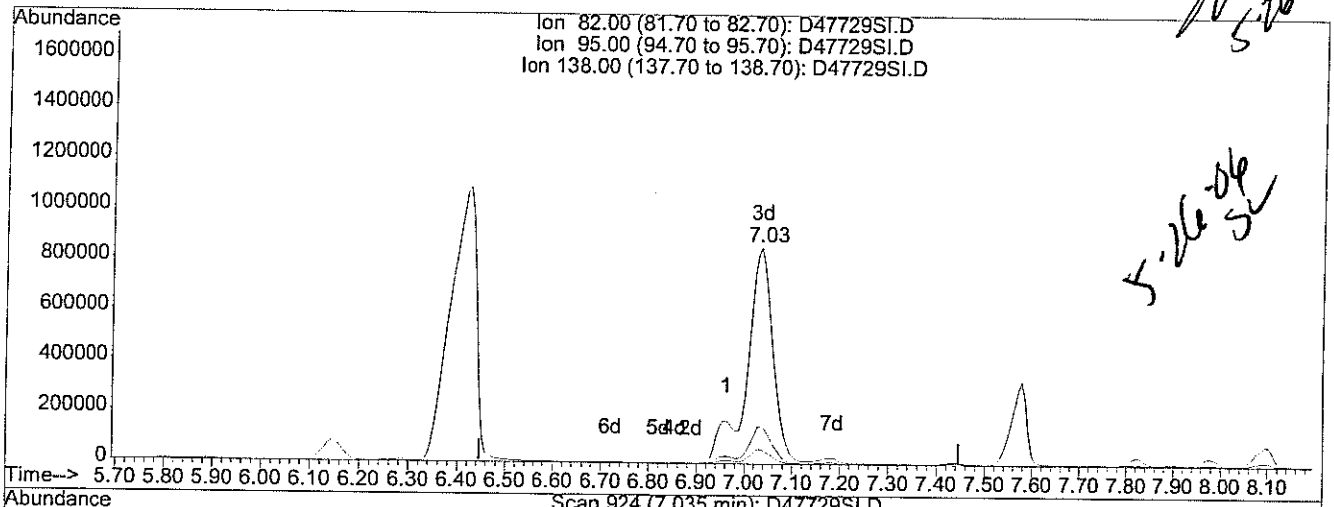
Data File : D:\HPCHEM\1\DATA\052406-D\D47729SI.D
Acq On : 25 May 2006 7:06 am
Sample : S6846 PT5 100/200
Misc :
MS Integration Params: RTEINT.P
Quant Time: May 26 9:27 2006

Vial: 8
Operator:
Inst : D Instrum
Multiplr: 1.00

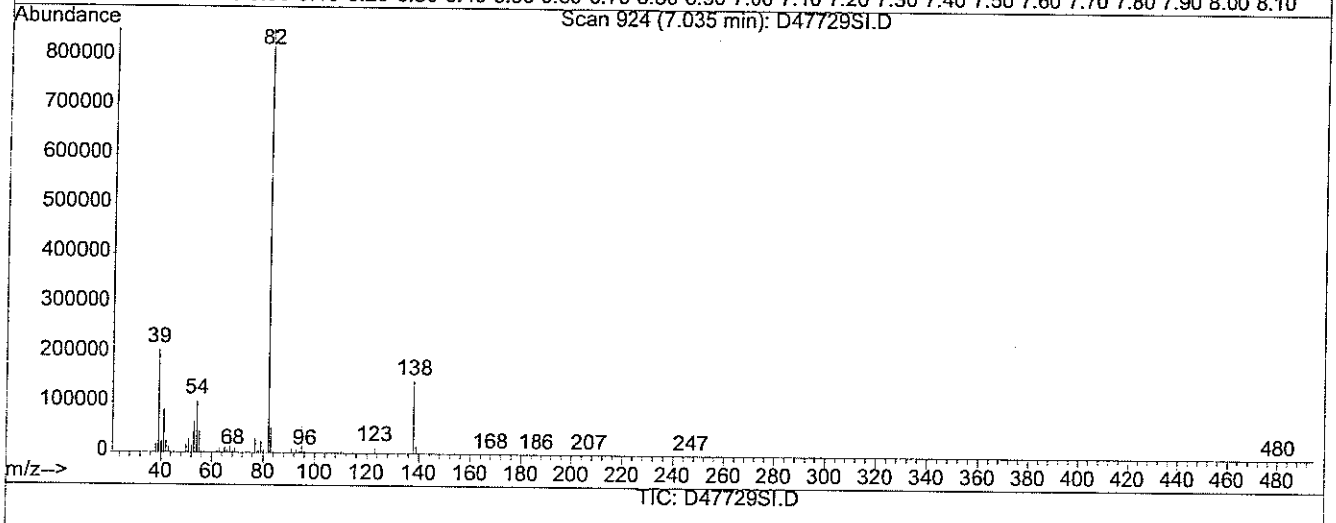
Quant Results File: temp.res

Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Fri May 26 09:21:45 2006
Response via : Multiple Level Calibration

5-26-06



5-26-06 SL



TIC: D47729SI.D

(22) Isophorone (TM)

7.03min 91.67ug/ml m

response 3325385

Ion	Exp%	Act%
82.00	100	100
95.00	6.20	1.05#
138.00	16.50	2.25#
0.00	0.00	0.00

Quantitation Report (Qedit)

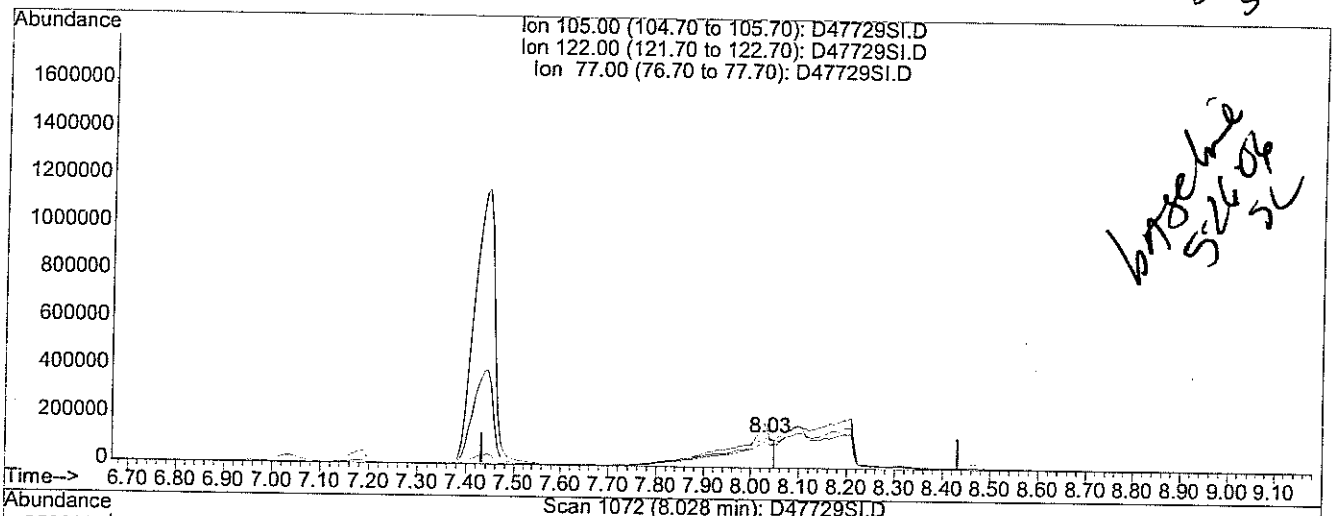
Data File : D:\HPCHEM\1\DATA\052406-D\D47729SI.D
Acq On : 25 May 2006 7:06 am
Sample : S6846 PT5 100/200
Misc :
MS Integration Params: RTEINT.P
Quant Time: May 26 9:27 2006

Vial: 8
Operator:
Inst : D Instrum
Multiplr: 1.00

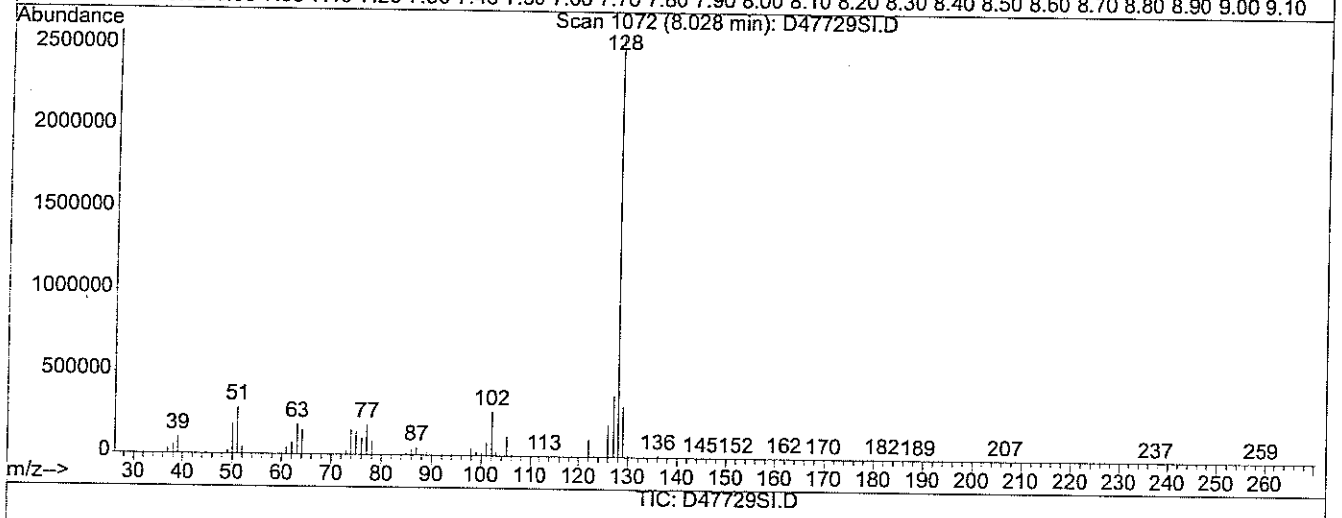
Quant Results File: temp.res

Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Fri May 26 09:21:45 2006
Response via : Multiple Level Calibration

5-26-06



baseline 5-26-06 SL



(27) Benzoic Acid (TM)

8.03min 90.66ug/ml

response 1066967

Ion	Exp%	Act%
105.00	100	100
122.00	78.50	74.61
77.00	70.70	55.06#
0.00	0.00	0.00

Quantitation Report (Qedit)

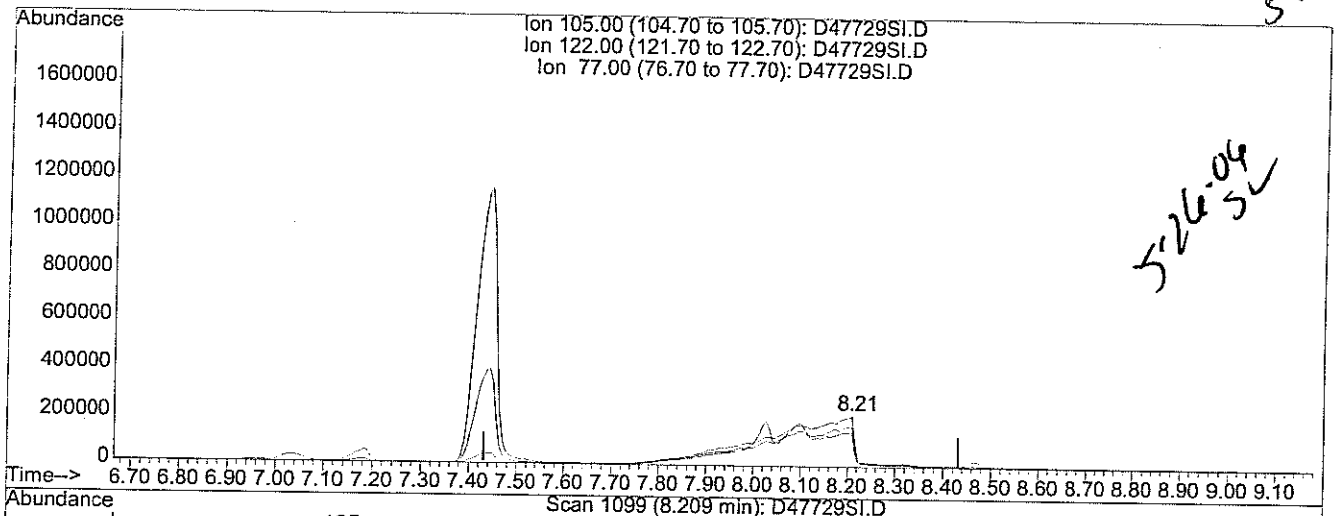
Data File : D:\HPCHEM\1\DATA\052406-D\D47729SI.D
Acq On : 25 May 2006 7:06 am
Sample : S6846 PT5 100/200
Misc :
MS Integration Params: RTEINT.P
Quant Time: May 26 9:27 2006

Vial: 8
Operator:
Inst : D Instrum
Multiplr: 1.00

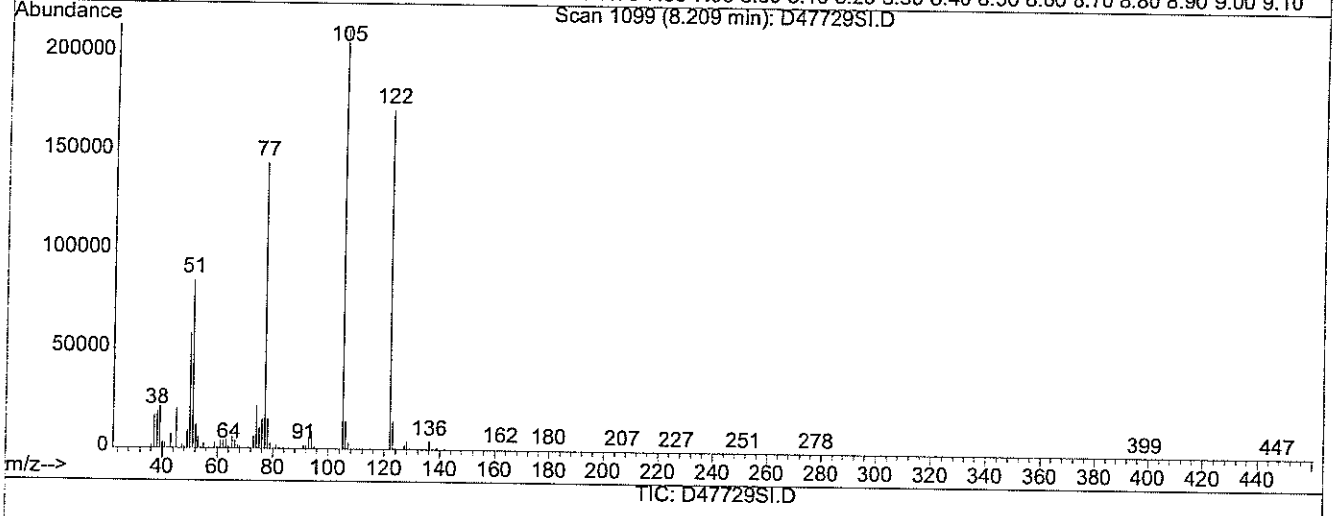
Quant Results File: temp.res

Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Fri May 26 09:21:45 2006
Response via : Multiple Level Calibration

5.26.06



5.26.06 SV



(27) Benzoic Acid (TM)

8.21min 223.77ug/ml m

response 2859129

Ion	Exp%	Act%
105.00	100	100
122.00	78.50	27.84#
77.00	70.70	20.55#
0.00	0.00	0.00

Quantitation Report (Qedit)

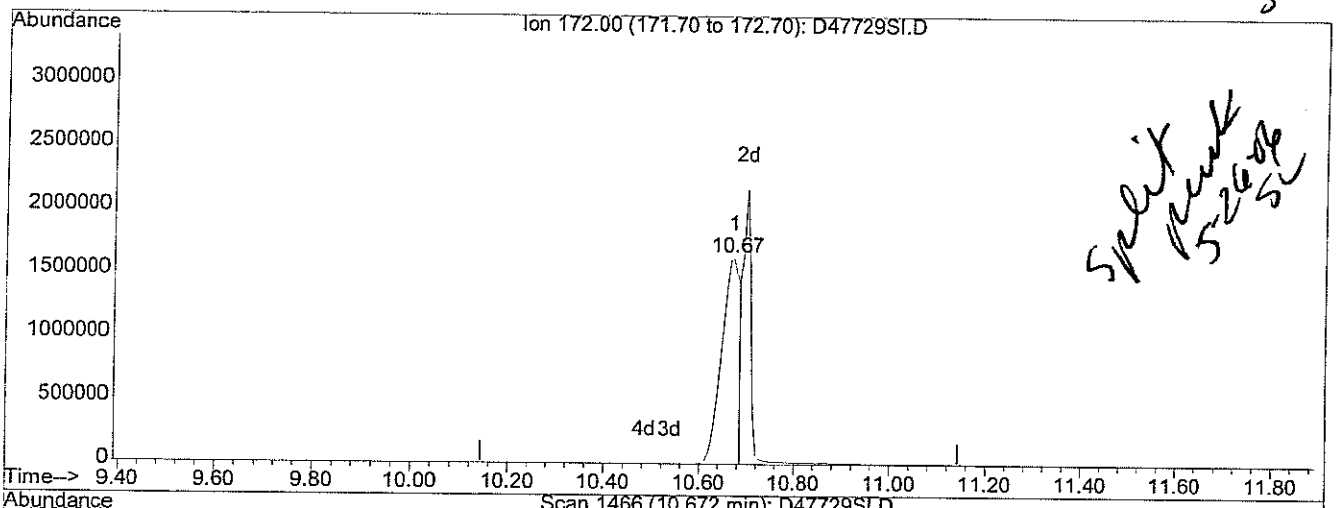
Data File : D:\HPCHEM\1\DATA\052406-D\D47729SI.D
Acq On : 25 May 2006 7:06 am
Sample : S6846 PT5 100/200
Misc :
MS Integration Params: RTEINT.P
Quant Time: May 26 9:27 2006

Vial: 8
Operator:
Inst : D Instrum
Multiplr: 1.00

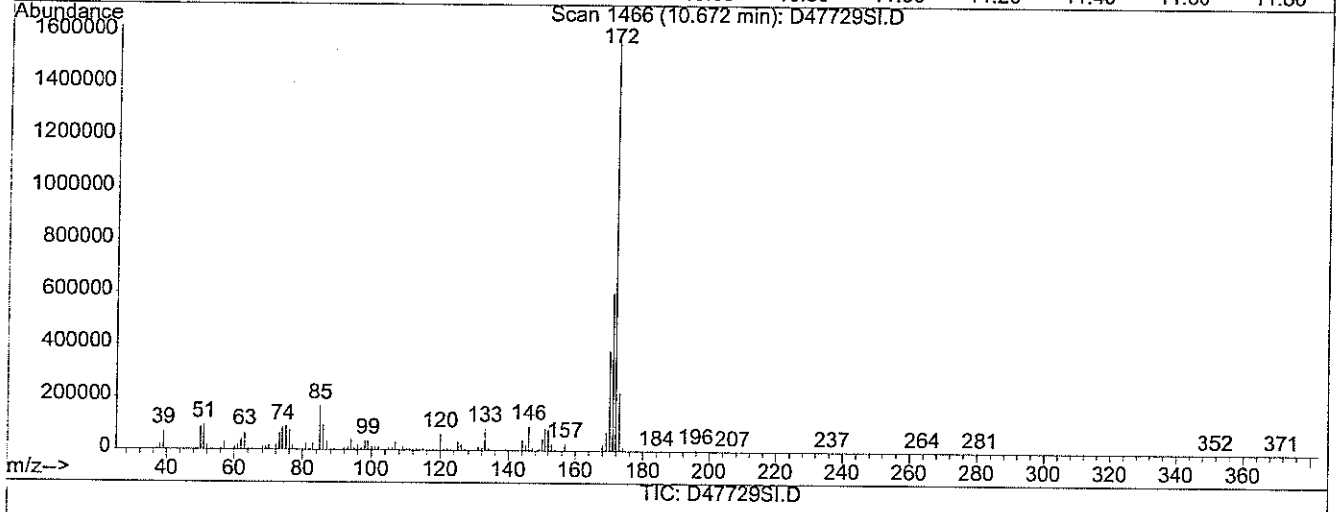
Quant Results File: temp.res

Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Fri May 26 09:21:45 2006
Response via : Multiple Level Calibration

5-26-06



Split Peak 5-26-06



TIC: D47729SI.D

(39) 2-Fluorobiphenyl (S)

10.67min 138.65ug/ml

response 4321236

Ion	Exp%	Act%
172.00	100	100
0.00	0.00	0.00
0.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (Qedit)

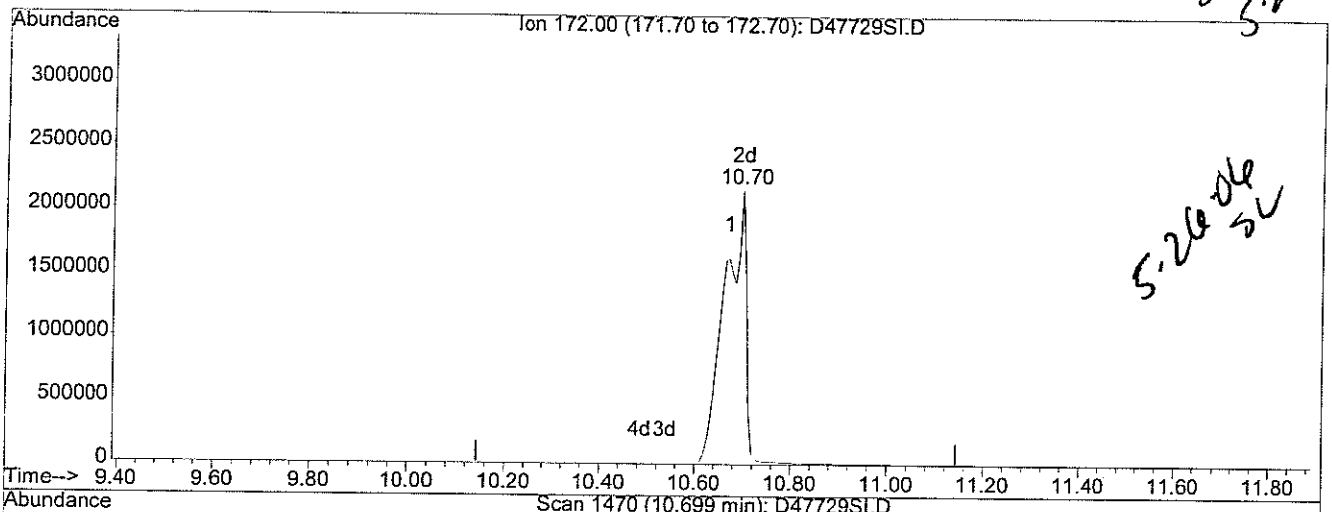
Data File : D:\HPCHEM\1\DATA\052406-D\D47729SI.D
Acq On : 25 May 2006 7:06 am
Sample : S6846 PT5 100/200
Misc :
MS Integration Params: RTEINT.P
Quant Time: May 26 9:27 2006

Vial: 8
Operator:
Inst : D Instrum
Multiplr: 1.00

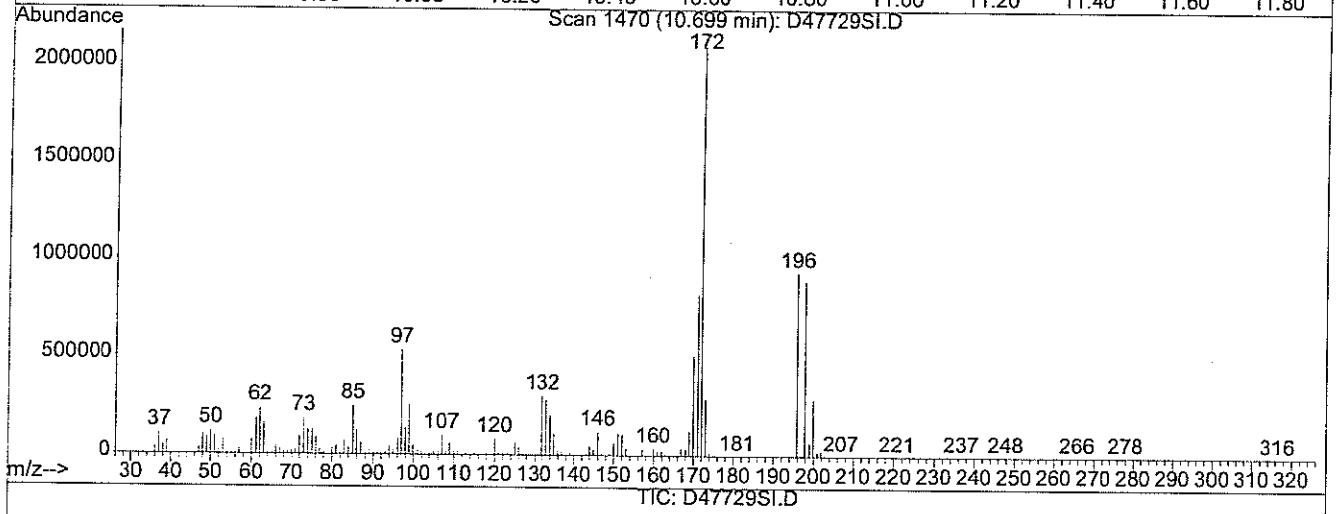
Quant Results File: temp.res

Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Fri May 26 09:21:45 2006
Response via : Multiple Level Calibration

5.26.06



5.26.06 SL



TIC: D47729SI.D

(39) 2-Fluorobiphenyl (S)

10.70min 218.27ug/ml m

response 6802770

Ion	Exp%	Act%
172.00	100	100
0.00	0.00	0.00
0.00	0.00	0.00
0.00	0.00	0.00

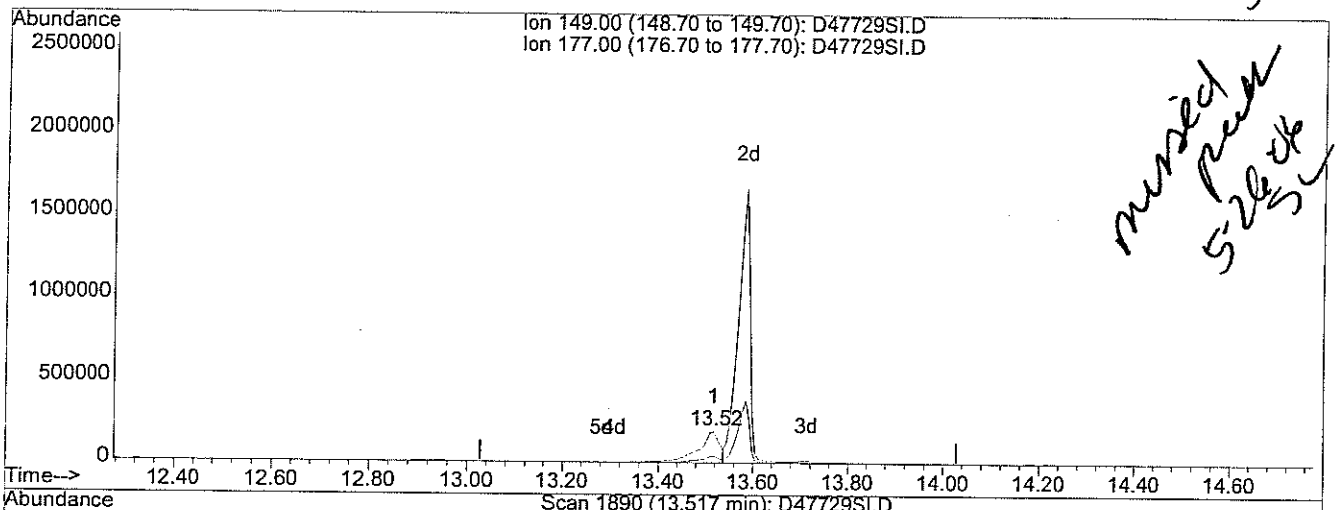
Quantitation Report (Qedit)

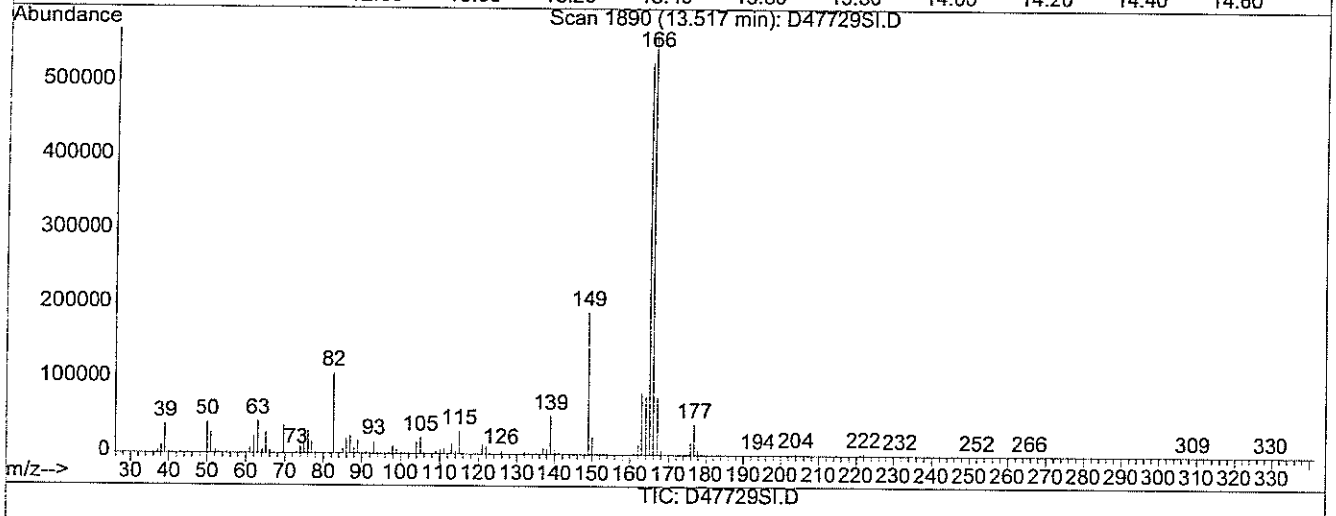
Data File : D:\HPCHEM\1\DATA\052406-D\D47729SI.D
Acq On : 25 May 2006 7:06 am
Sample : S6846 PT5 100/200
Misc :
MS Integration Params: RTEINT.P
Quant Time: May 26 9:27 2006

Vial: 8
Operator:
Inst : D Instrum
Multiplr: 1.00

Quant Results File: temp.res

Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Fri May 26 09:21:45 2006
Response via : Multiple Level Calibration





TIC: D47729SI.D

(53) Diethyl Phthalate (TM)

13.52min 16.78ug/ml

response 606206

Ion	Exp%	Act%
149.00	100	100
177.00	21.70	21.26
0.00	0.00	0.00
0.00	0.00	0.00

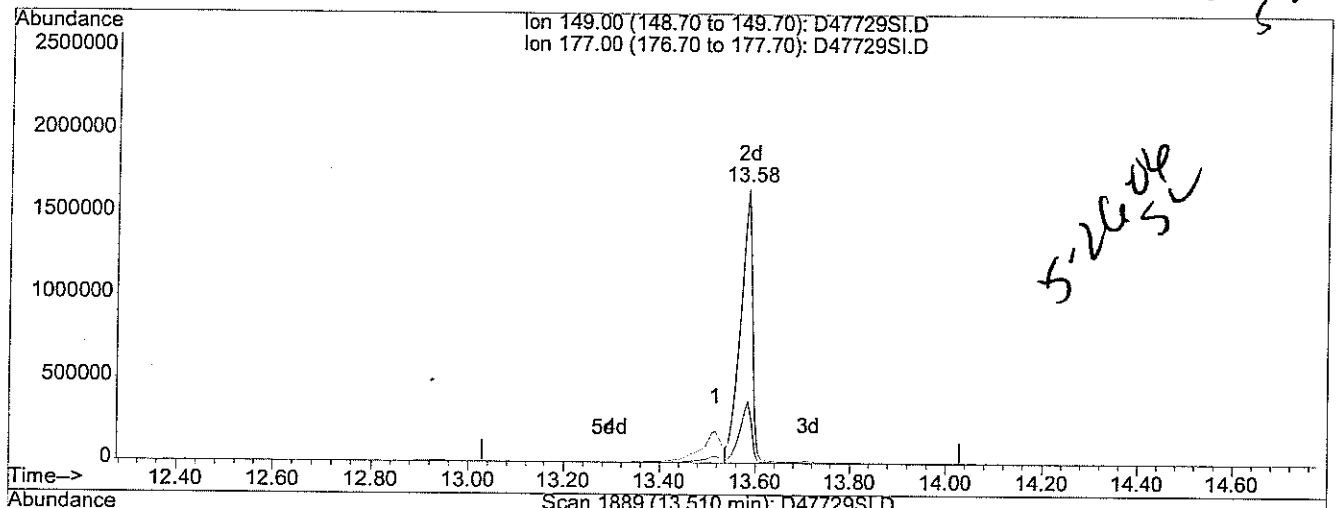
Quantitation Report (Qedit)

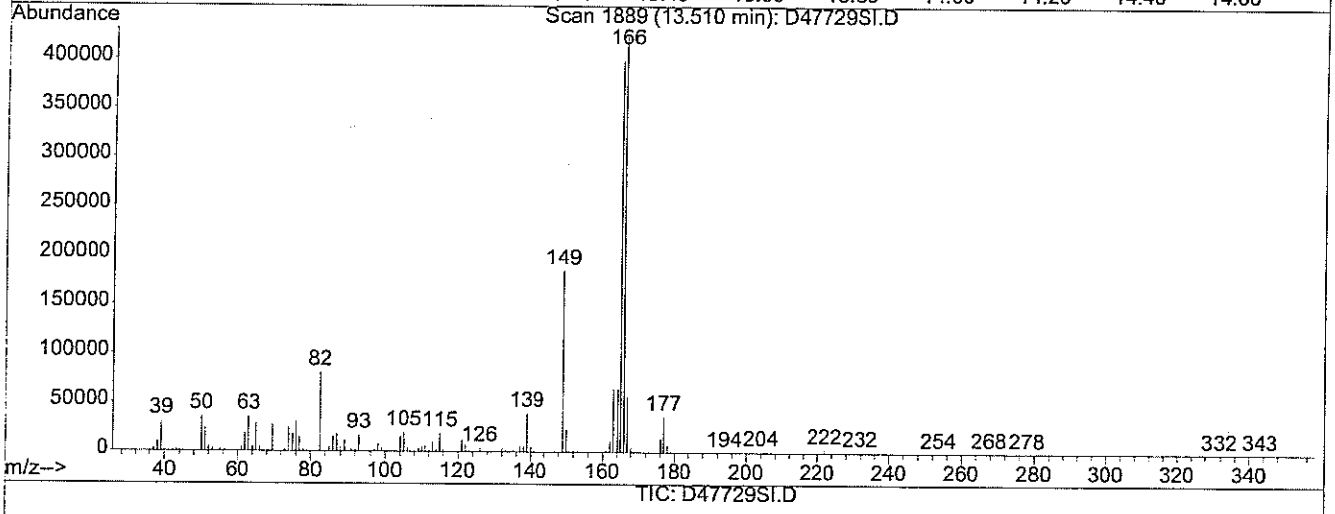
Data File : D:\HPCHEM\1\DATA\052406-D\D47729SI.D
Acq On : 25 May 2006 7:06 am
Sample : S6846 PT5 100/200
Misc :
MS Integration Params: RTEINT.P
Quant Time: May 26 9:28 2006

Vial: 8
Operator:
Inst : D Instrum
Multiplr: 1.00

Quant Results File: temp.res

Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Fri May 26 09:21:45 2006
Response via : Multiple Level Calibration





TIC: D47729SI.D

(53) Diethyl Phthalate (TM)

13.58min 79.65ug/ml m

response 2877020

Ion	Exp%	Act%
149.00	100	100
177.00	21.70	4.48#
0.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (Qedit)

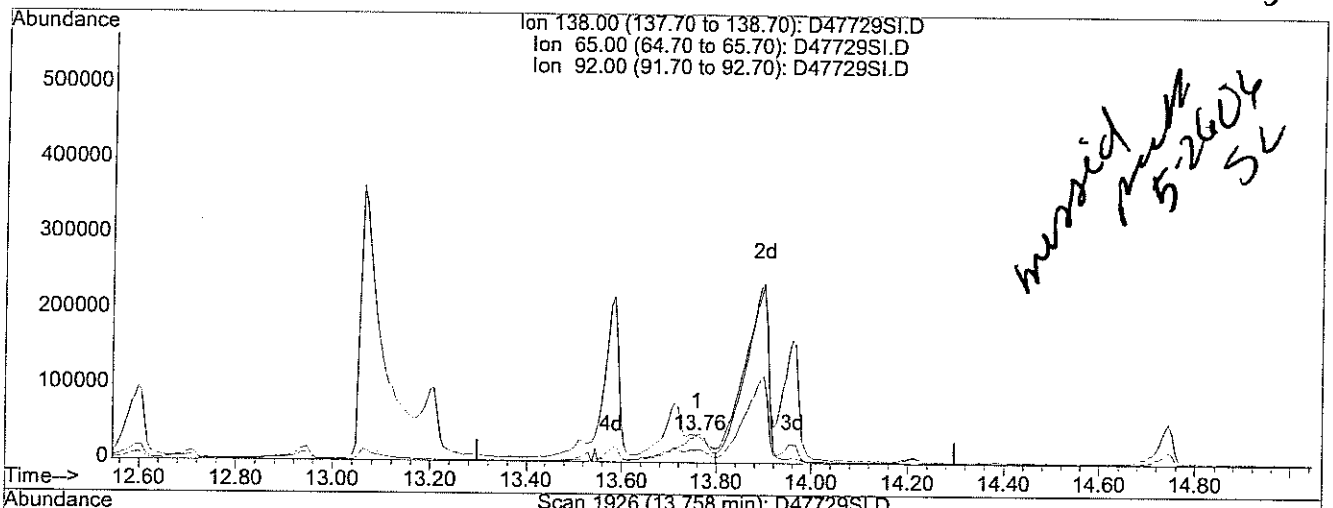
Data File : D:\HPCHEM\1\DATA\052406-D\D47729SI.D
Acq On : 25 May 2006 7:06 am
Sample : S6846 PT5 100/200
Misc :
MS Integration Params: RTEINT.P
Quant Time: May 26 9:28 2006

Vial: 8
Operator:
Inst : D Instrum
Multiplr: 1.00

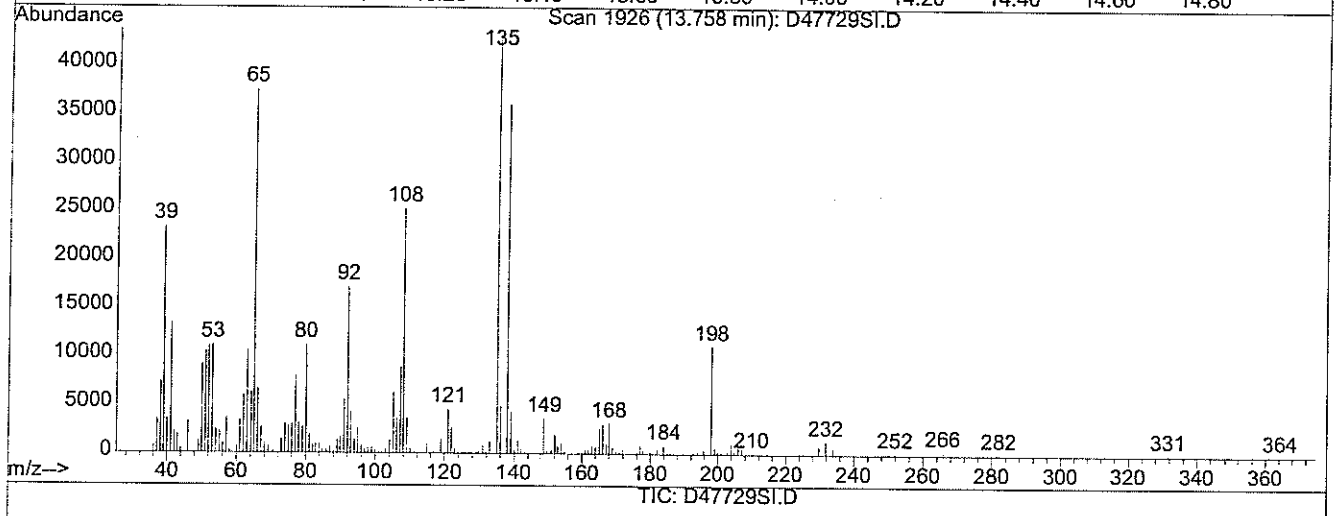
Quant Results File: temp.res

Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Fri May 26 09:21:45 2006
Response via : Multiple Level Calibration

5.26.06



Incorrect peak 5.26.06 SL



TIC: D47729SI.D

(55) 4-Nitroaniline (TM)

13.76min 17.61ug/ml

response 197602

Ion	Exp%	Act%
138.00	100	100
65.00	96.40	94.74
92.00	47.80	27.16#
0.00	0.00	0.00

Quantitation Report (Qedit)

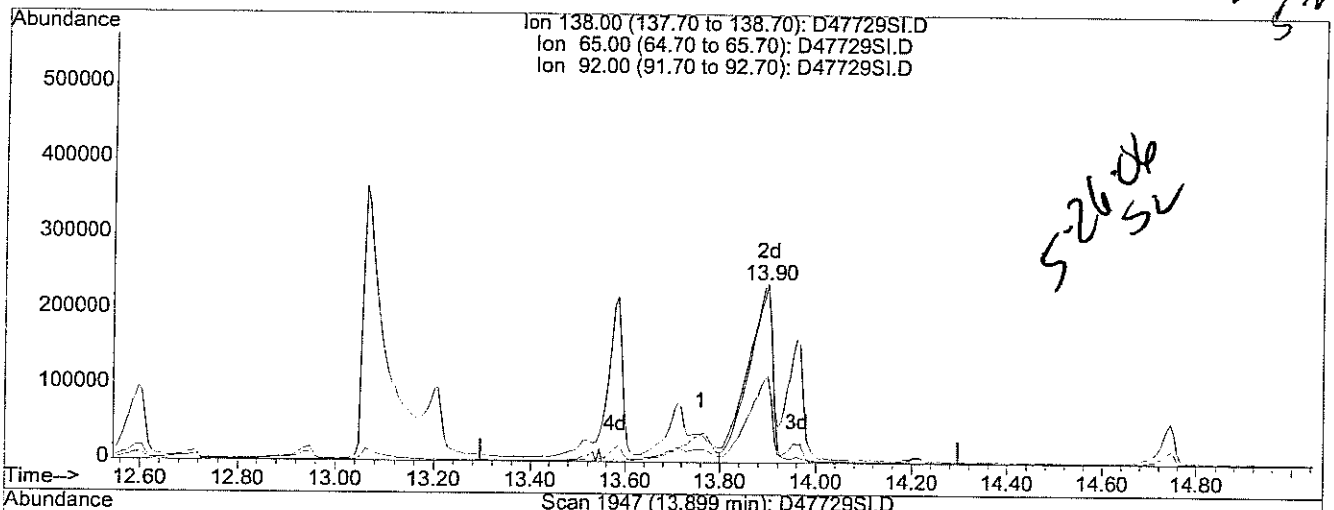
Data File : D:\HPCHEM\1\DATA\052406-D\D47729SI.D
Acq On : 25 May 2006 7:06 am
Sample : S6846 PT5 100/200
Misc :
MS Integration Params: RTEINT.P
Quant Time: May 26 9:28 2006

Vial: 8
Operator:
Inst : D Instrum
Multiplr: 1.00

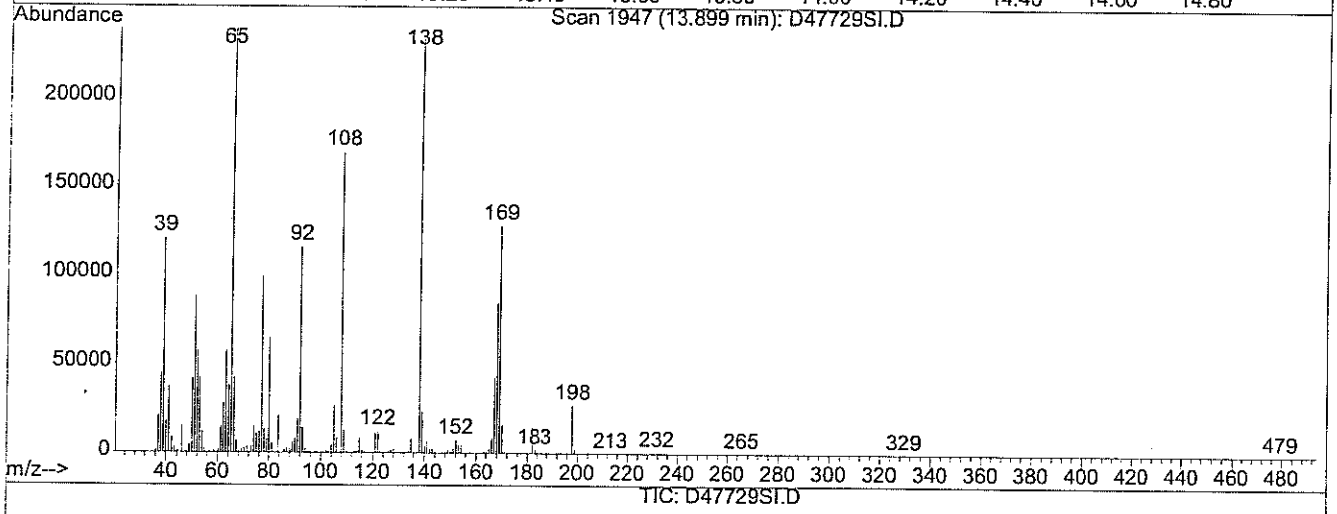
Quant Results File: temp.res

Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Fri May 26 09:21:45 2006
Response via : Multiple Level Calibration

5-26-06



5-26-06 SL



(55) 4-Nitroaniline (TM)

13.90min 73.32ug/ml m

response 822653

Ion	Exp%	Act%
138.00	100	100
65.00	96.40	22.76#
92.00	47.80	6.52#
0.00	0.00	0.00

Quantitation Report (Qedit)

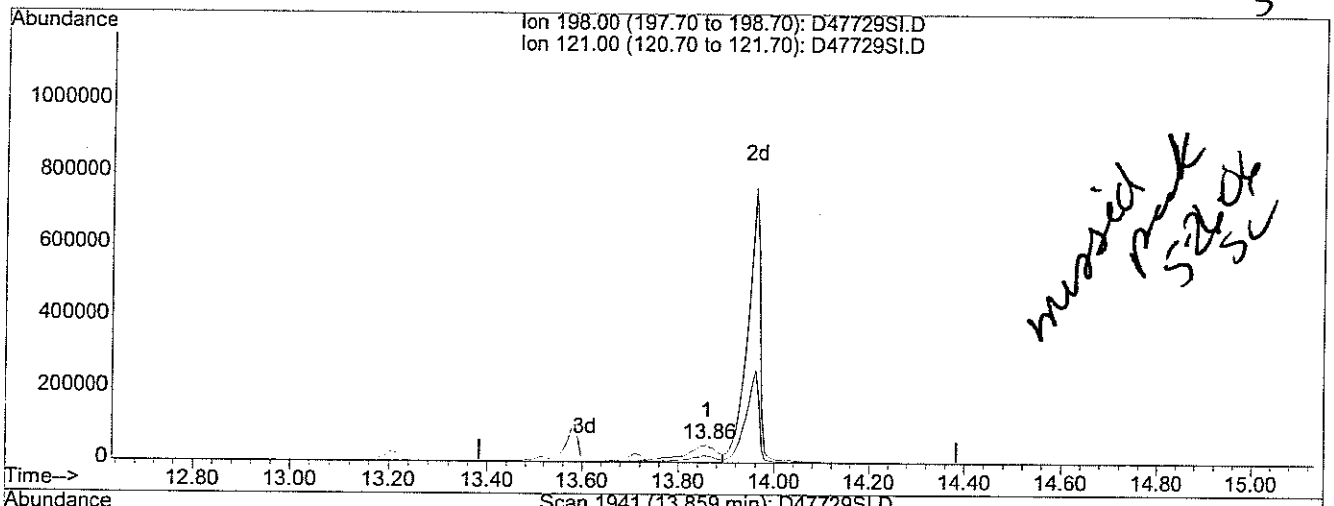
Data File : D:\HPCHEM\1\DATA\052406-D\D47729SI.D
Acq On : 25 May 2006 7:06 am
Sample : S6846 PT5 100/200
Misc :
MS Integration Params: RTEINT.P
Quant Time: May 26 9:28 2006

Vial: 8
Operator:
Inst : D Instrum
Multiplr: 1.00

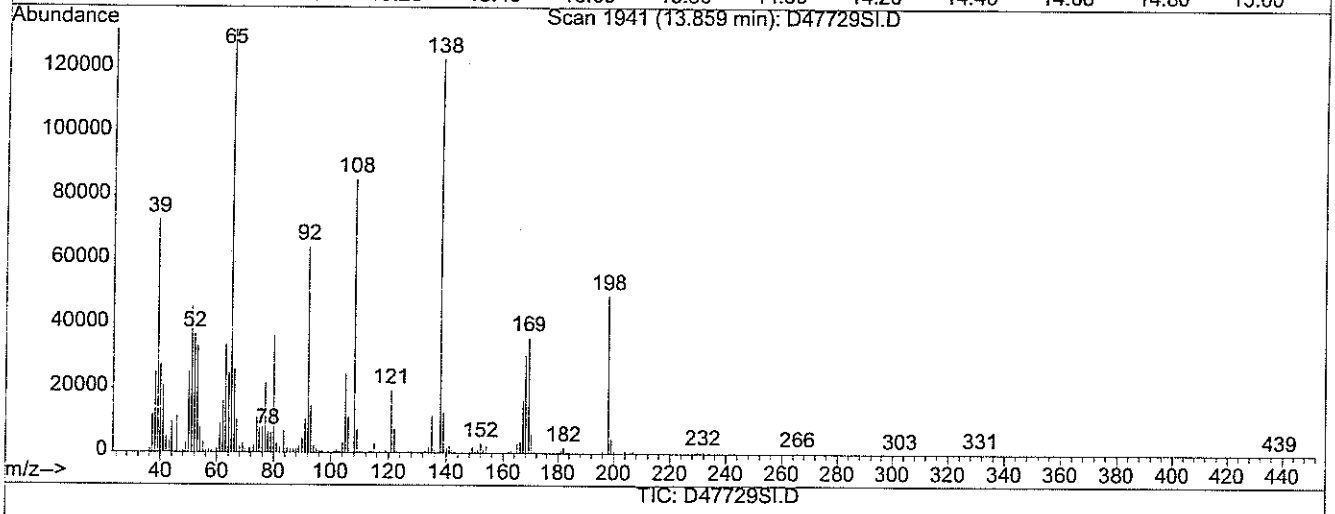
Quant Results File: temp.res

Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Fri May 26 09:21:45 2006
Response via : Multiple Level Calibration

5-26-06



missed peak 5-26-06



TIC: D47729SI.D

(56) 4,6-Dinitro-2-methylphenol (TM)

13.86min 30.13ug/ml

response 236312

Ion	Exp%	Act%
198.00	100	100
121.00	37.30	31.21
0.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (Qedit)

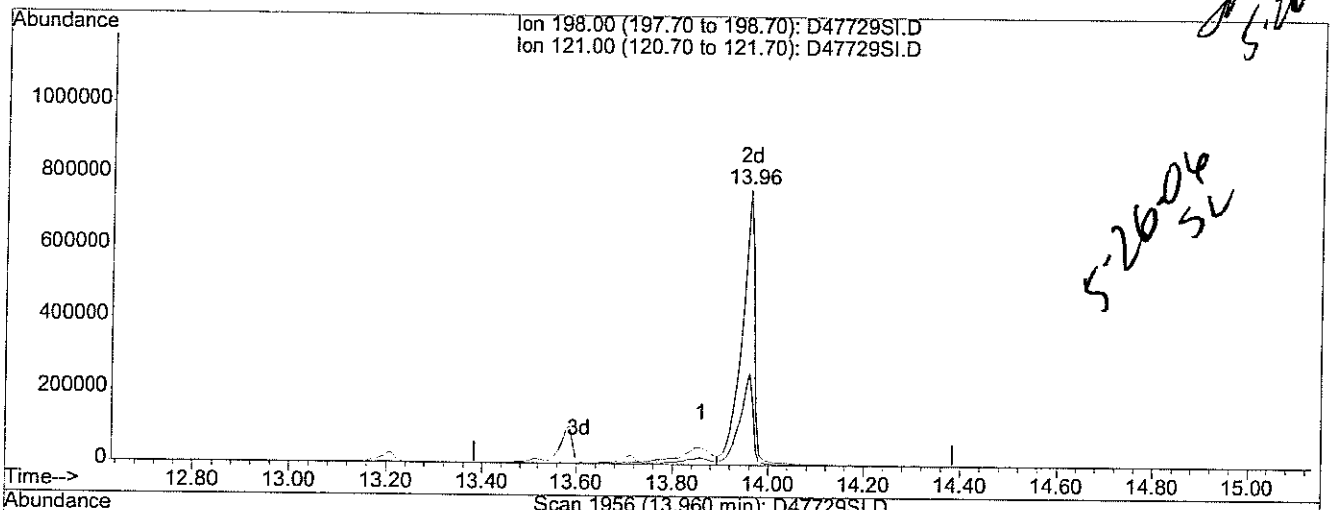
Data File : D:\HPCHEM\1\DATA\052406-D\D47729SI.D
Acq On : 25 May 2006 7:06 am
Sample : S6846 PT5 100/200
Misc :
MS Integration Params: RTEINT.P
Quant Time: May 26 9:28 2006

Vial: 8
Operator:
Inst : D Instrum
Multiplr: 1.00

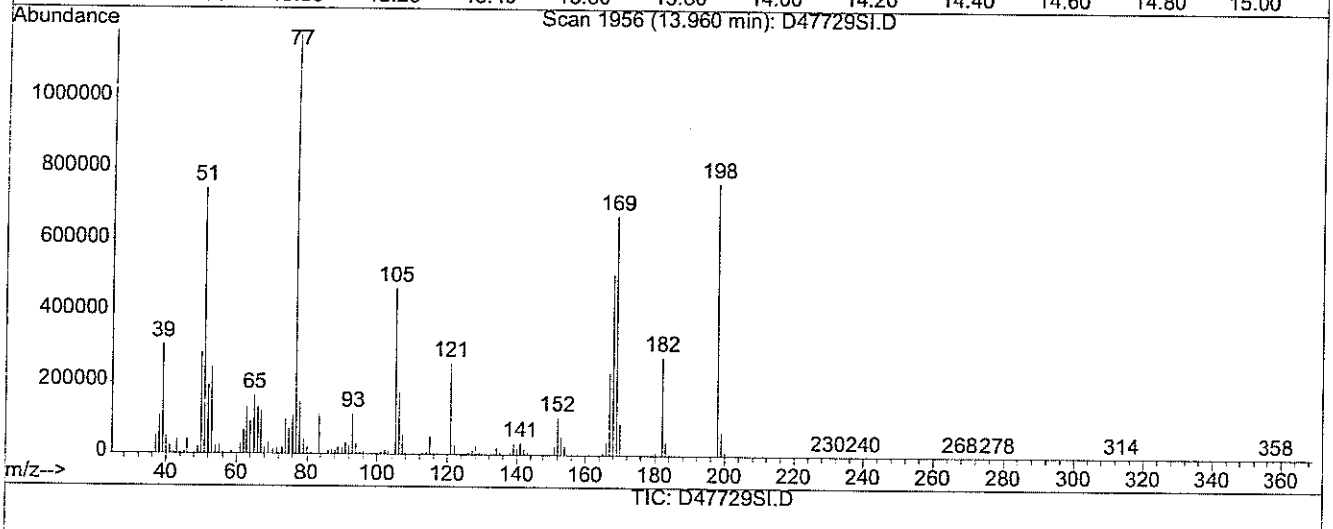
Quant Results File: temp.res

Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Fri May 26 09:21:45 2006
Response via : Multiple Level Calibration

5-26-06



5-26-06 SL



TIC: D47729SI.D

(56) 4,6-Dinitro-2-methylphenol (TM)

13.96min 219.23ug/ml m

response 1449894

Ion	Exp%	Act%
198.00	100	100
121.00	37.30	5.09#
0.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (Qedit)

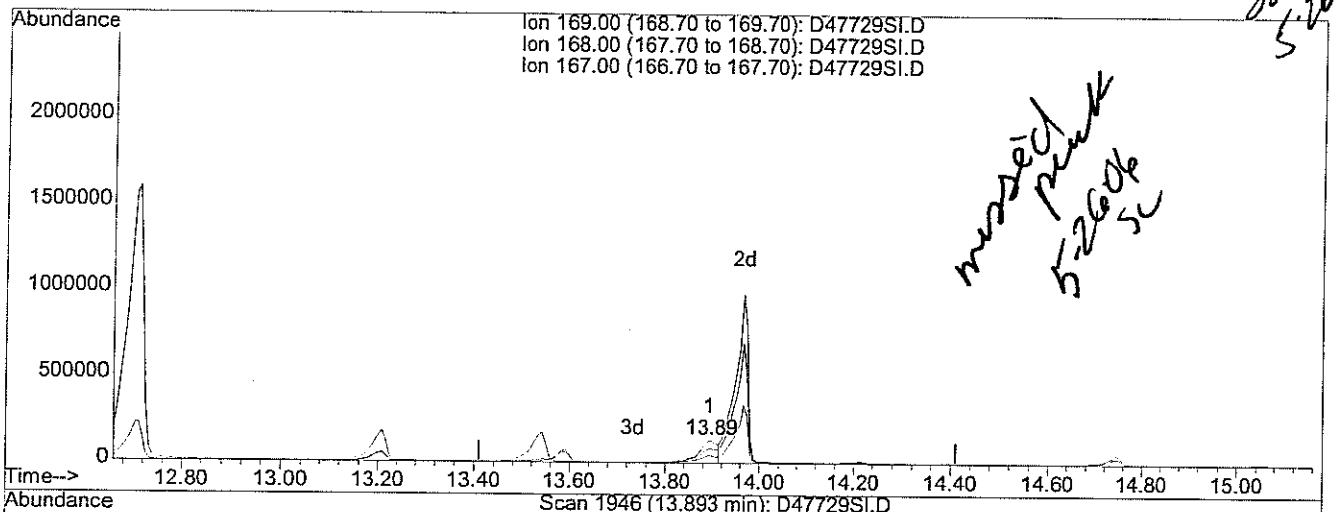
Data File : D:\HPCHEM\1\DATA\052406-D\D47729SI.D
Acq On : 25 May 2006 7:06 am
Sample : S6846 PT5 100/200
Misc :
MS Integration Params: RTEINT.P
Quant Time: May 26 9:28 2006

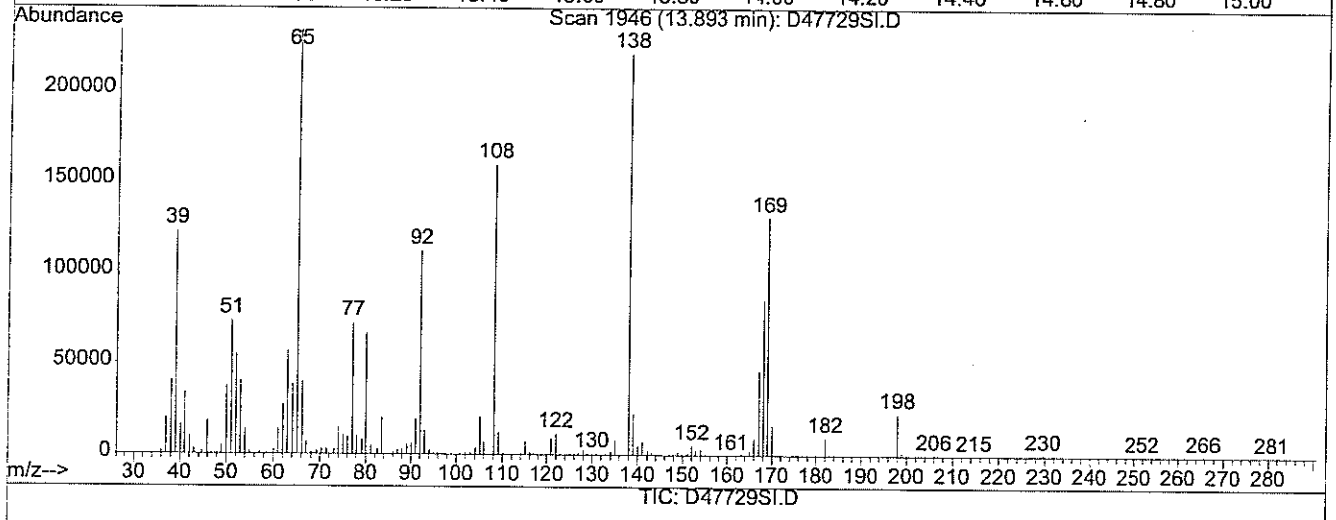
Vial: 8
Operator:
Inst : D Instrum
Multiplr: 1.00

Quant Results File: temp.res

Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Fri May 26 09:21:45 2006
Response via : Multiple Level Calibration

5.26.06





(57) n-nitrosodiphenylamine (TCM)

13.89min 17.66ug/ml

response 373269

Ion	Exp%	Act%
169.00	100	100
168.00	68.80	67.06
167.00	33.20	39.80
0.00	0.00	0.00

Quantitation Report (Qedit)

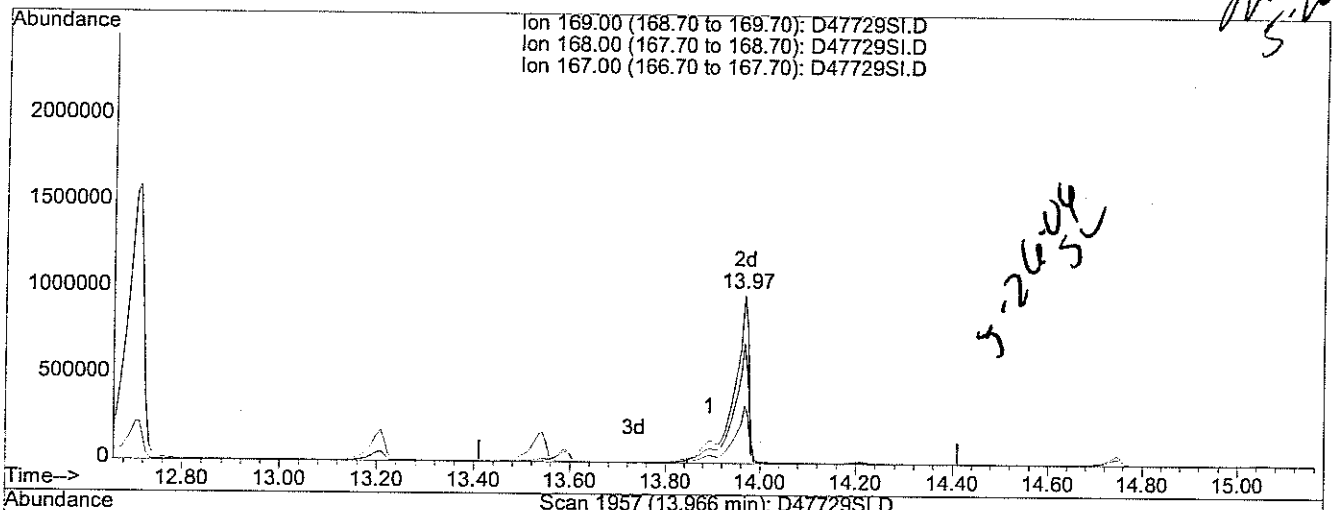
Data File : D:\HPCHEM\1\DATA\052406-D\D47729SI.D
Acq On : 25 May 2006 7:06 am
Sample : S6846 PT5 100/200
Misc :
MS Integration Params: RTEINT.P
Quant Time: May 26 9:28 2006

Vial: 8
Operator:
Inst : D Instrum
Multiplr: 1.00

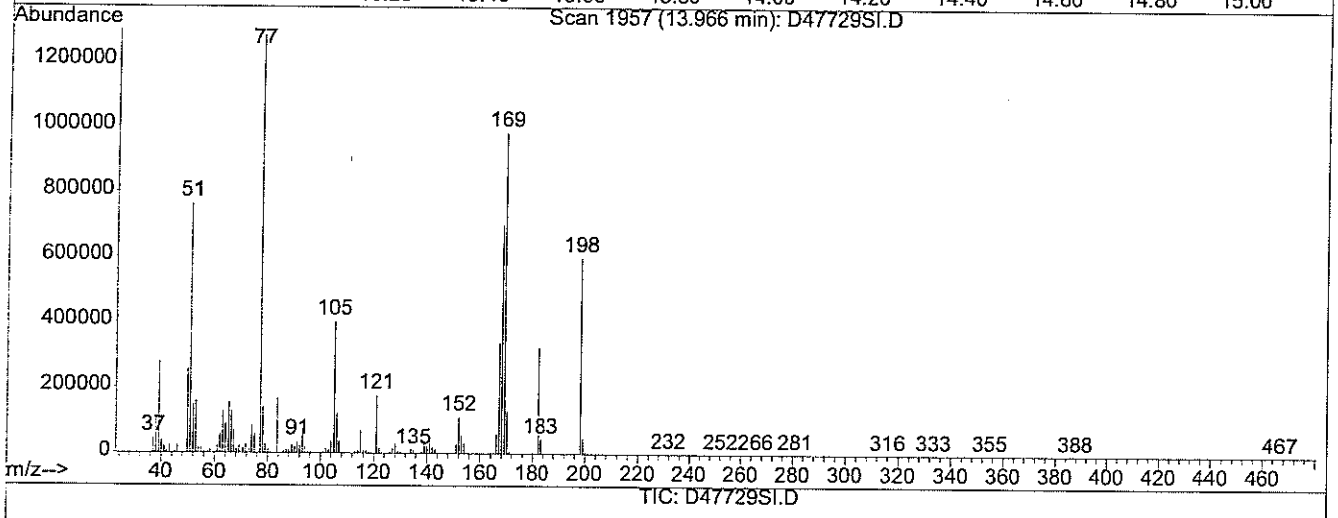
Quant Results File: temp.res

Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Fri May 26 09:21:45 2006
Response via : Multiple Level Calibration

5-26-06



5-26-06



TIC: D47729SI.D

(57) n-nitrosodiphenylamine (TCM)

13.97min 105.08ug/ml m

response 2220590

Ion	Exp%	Act%
169.00	100	100
168.00	68.80	11.27#
167.00	33.20	6.69#
0.00	0.00	0.00

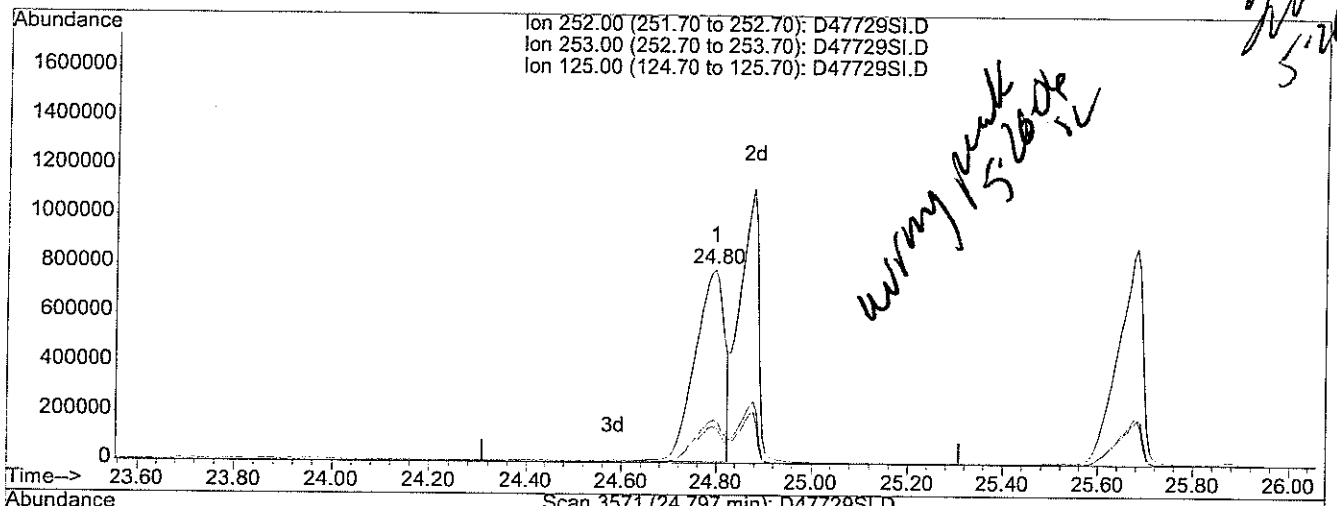
Quantitation Report (Qedit)

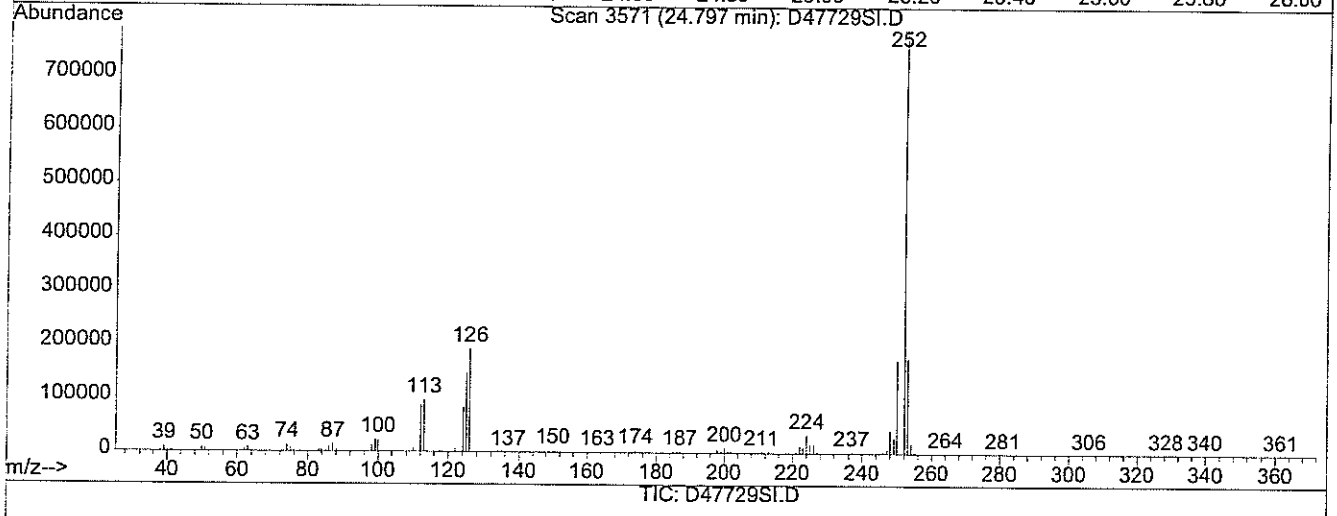
Data File : D:\HPCHEM\1\DATA\052406-D\D47729SI.D
Acq On : 25 May 2006 7:06 am
Sample : S6846 PT5 100/200
Misc :
MS Integration Params: RTEINT.P
Quant Time: May 26 9:28 2006

Vial: 8
Operator:
Inst : D Instrum
Multiplr: 1.00

Quant Results File: temp.res

Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Fri May 26 09:21:45 2006
Response via : Multiple Level Calibration





TIC: D47729SI.D

(82) Benzo[k] fluoranthene (TM)

24.80min 98.23ug/ml

response 3368715

Ion	Exp%	Act%
252.00	100	100
253.00	22.20	22.11
125.00	18.50	18.81
0.00	0.00	0.00

Quantitation Report (Qedit)

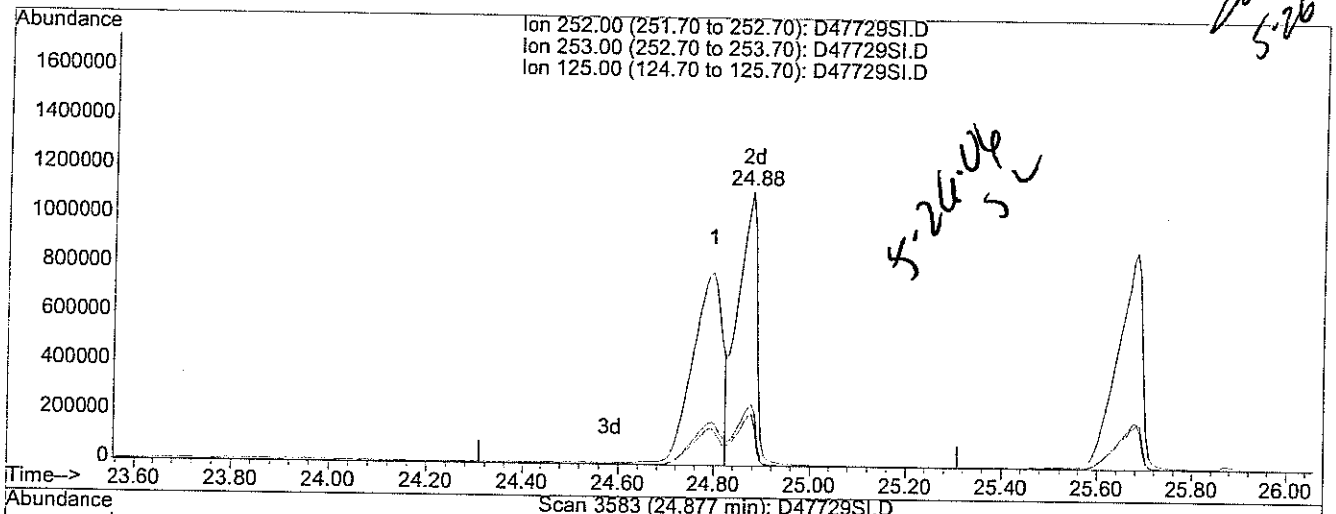
Data File : D:\HPCHEM\1\DATA\052406-D\D47729SI.D
Acq On : 25 May 2006 7:06 am
Sample : S6846 PT5 100/200
Misc :
MS Integration Params: RTEINT.P
Quant Time: May 26 9:28 2006

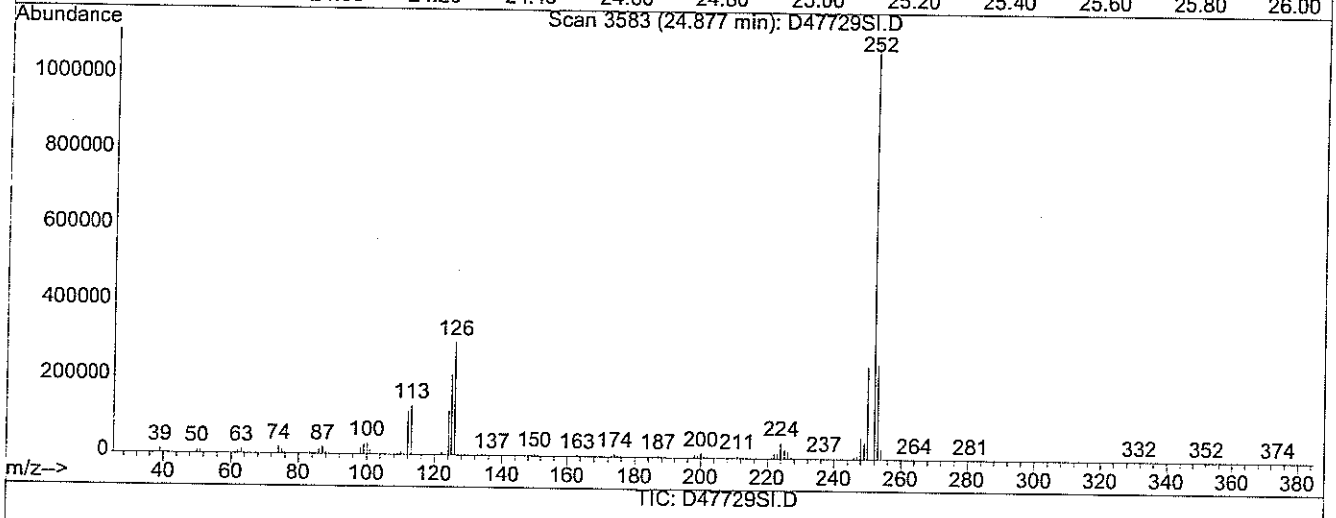
Vial: 8
Operator:
Inst : D Instrum
Multiplr: 1.00

Quant Results File: temp.res

Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Fri May 26 09:21:45 2006
Response via : Multiple Level Calibration

5-26-06





(82) Benzo[k] fluoranthene (TM)

24.88min 93.14ug/ml m

response 3194200

Ion	Exp%	Act%
252.00	100	100
253.00	22.20	23.32
125.00	18.50	19.84
0.00	0.00	0.00

Quantitation Report (Qedit)

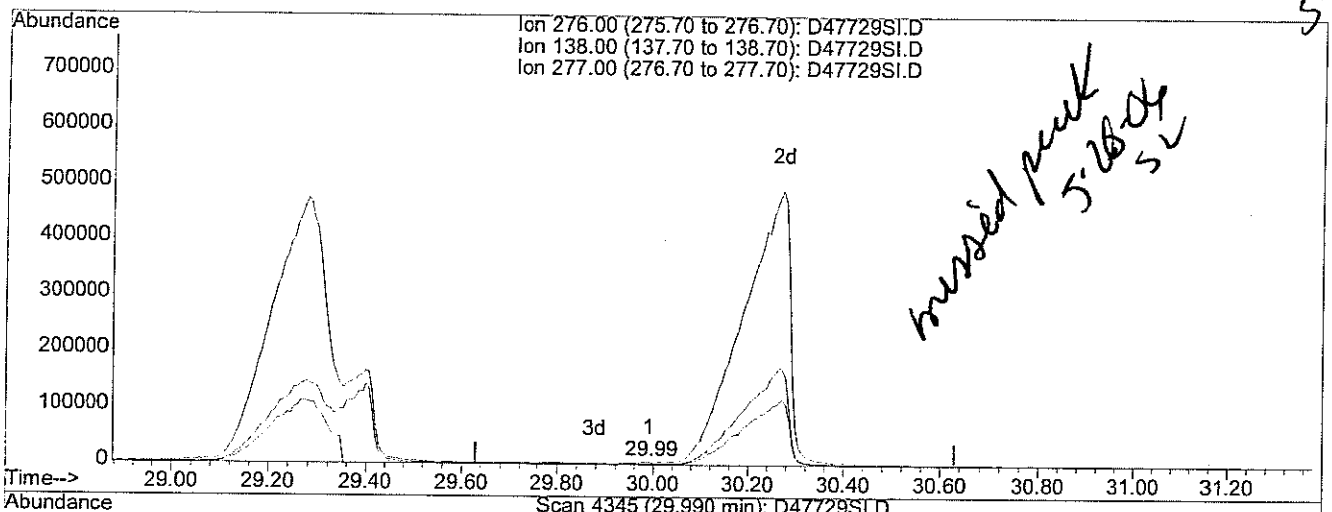
Data File : D:\HPCHEM\1\DATA\052406-D\D47729SI.D
Acq On : 25 May 2006 7:06 am
Sample : S6846 PT5 100/200
Misc :
MS Integration Params: RTEINT.P
Quant Time: May 26 9:28 2006

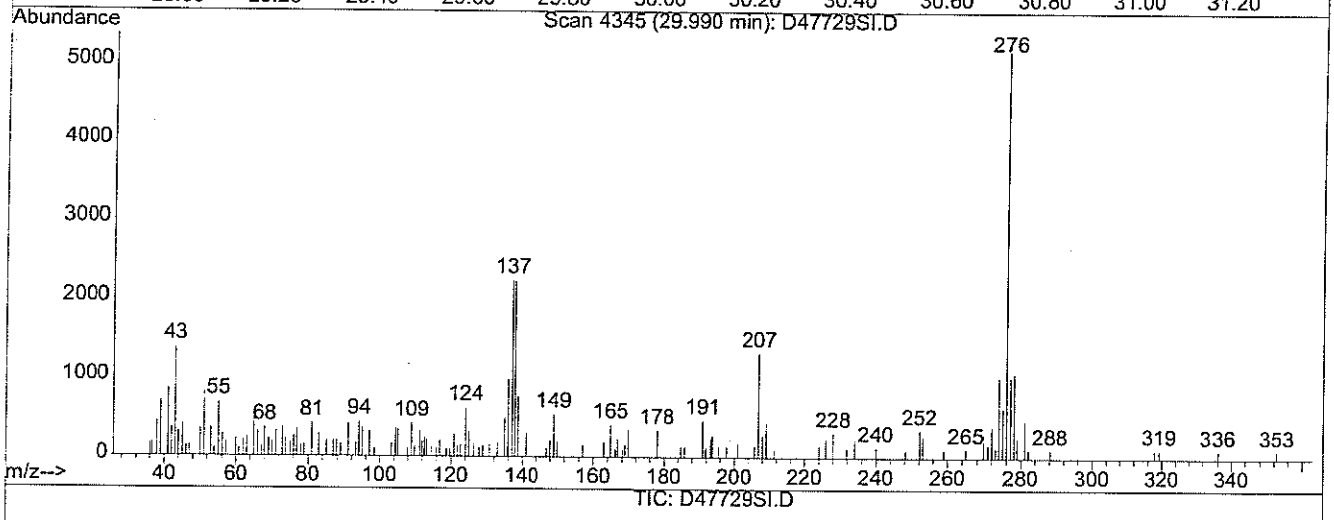
Vial: 8
Operator:
Inst : D Instrum
Multiplr: 1.00

Quant Results File: temp.res

Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Fri May 26 09:21:45 2006
Response via : Multiple Level Calibration

5-26-06





TIC: D47729SI.D

(86) Benzo(g,h,i) perylene (TM)

29.99min 0.26ug/ml

response 8935

Ion	Exp%	Act%
276.00	100	100
138.00	0.10	19.22#
277.00	0.00	21.97#
0.00	0.00	0.00

Quantitation Report (Qedit)

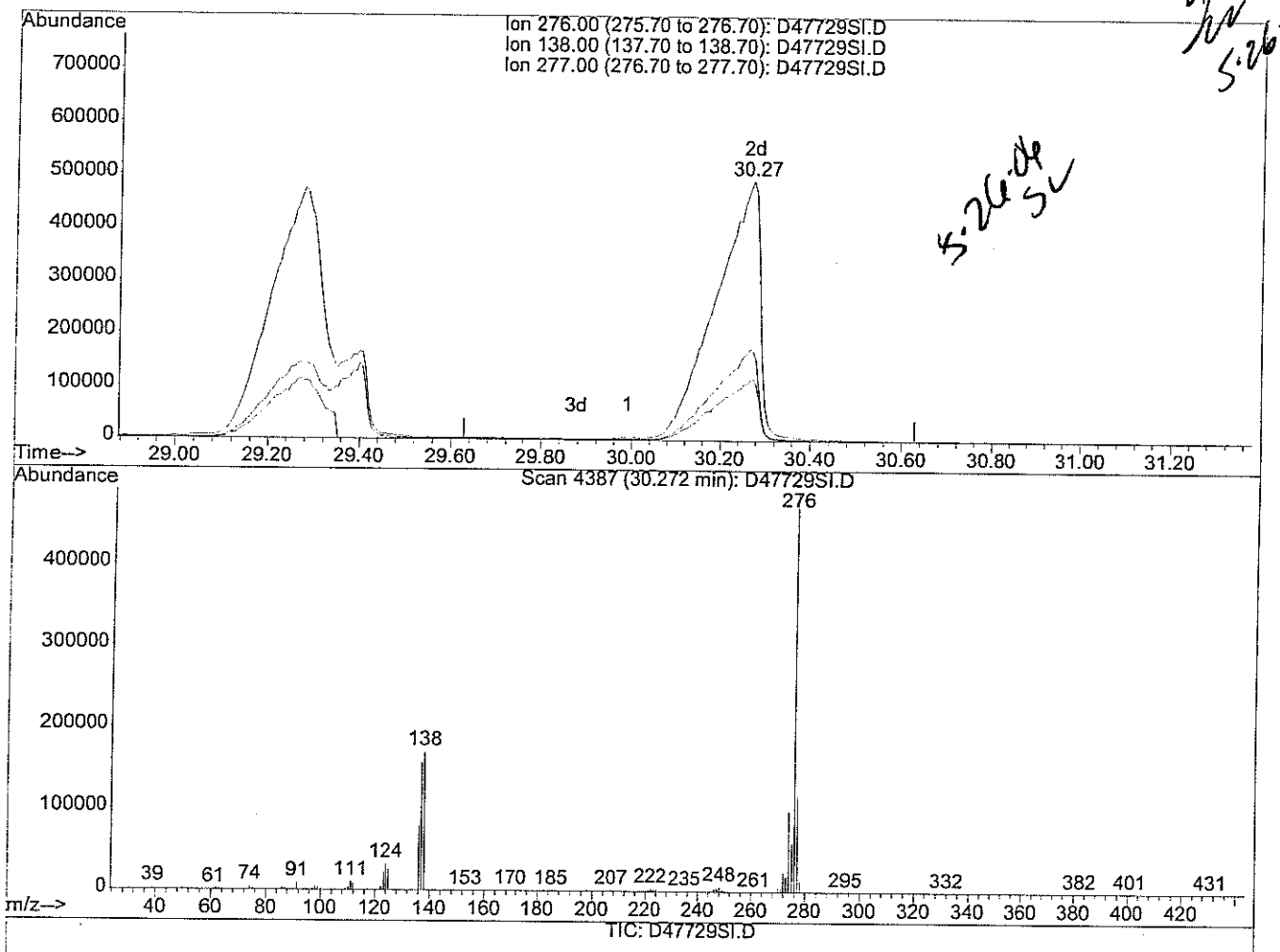
Data File : D:\HPCHEM\1\DATA\052406-D\D47729SI.D
Acq On : 25 May 2006 7:06 am
Sample : S6846 PT5 100/200
Misc :
MS Integration Params: RTEINT.P
Quant Time: May 26 9:29 2006

Vial: 8
Operator:
Inst : D Instrum
Multiplr: 1.00

Quant Results File: temp.res

Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Fri May 26 09:21:45 2006
Response via : Multiple Level Calibration

5.26.06



(86) Benzo(g,h,i) perylene (TM)

30.27min 100.65ug/ml m

response 3406455

Ion	Exp%	Act%
276.00	100	100
138.00	0.10	0.05#
277.00	0.00	0.06#
0.00	0.00	0.00

ABN CC

Data File Name D47730.D
Date Acquired 5/25/2006 7:39
Sample Name S6857 ABN CC 20/40

Name	ug/ml Amount	ug/ml Expected	% RECOVERY	70%-130%
Pyridine	19.60	20	98%	OK
N-nitrosodimethylamine	19.80	20	99%	OK
Phenol	40.30	40	101%	OK
2-Chlorophenol	40.96	40	102%	OK
bis(2-chloroethyl) ether	20.88	20	104%	OK
1,3-Dichlorobenzene	21.37	20	107%	OK
1,4-Dichlorobenzene	20.87	20	104%	OK
1,2-Dichlorobenzene	20.71	20	104%	OK
2-Methylphenol	40.78	40	102%	OK
bis(2-chloroisopropyl)ether	20.40	20	102%	OK
3+4-Methylphenol	42.15	40	105%	OK
Hexachloroethane	20.97	20	105%	OK
N-nitroso-di-n-propylamine	22.09	20	110%	OK
Nitrobenzene	21.18	20	106%	OK
Isophorone	23.56	20	118%	OK
2-Nitrophenol	42.99	40	107%	OK
2,4-Dimethylphenol	40.00	40	100%	OK
Bis(2-chloroethoxy)methane	20.48	20	102%	OK
2,4-Dichlorophenol	41.25	40	103%	OK
Benzoic Acid	39.94	40	100%	OK
1,2,4-Trichlorobenzene	20.90	20	105%	OK
Naphthalene	20.45	20	102%	OK
2,6-Dichlorophenol	40.34	40	101%	OK
Hexachlorobutadiene	21.26	20	106%	OK
4-Chloro-3-methylphenol	40.73	40	102%	OK
Hexachlorocyclopentadiene	25.14	20	126%	OK
2,4,6-Trichlorophenol	41.37	40	103%	OK
2,4,5-Trichlorophenol	38.77	40	97%	OK
2-Chloronaphthalene	20.37	20	102%	OK
2,6-Dinitrotoluene	21.09	20	105%	OK
Dimethyl Phthalate	20.53	20	103%	OK
Acenaphthylene	21.29	20	106%	OK
Acenaphthene	20.81	20	104%	OK
2,4-dinitrophenol	39.89	40	100%	OK
4-Nitrophenol	38.15	40	95%	OK
2,4-Dinitrotoluene	22.46	20	112%	OK
2,3,4,6-Tetrachlorophenol	42.19	40	105%	OK
Fluorene	20.78	20	104%	OK
Diethyl Phthalate	21.68	20	108%	OK
4-Chlorophenyl phenyl ether	20.73	20	104%	OK
4,6-Dinitro-2-methylphenol	39.80	40	100%	OK
n-nitrosodiphenylamine	20.09	20	100%	OK
Azobenzene	19.12	20	96%	OK
4-bromophenyl phenyl ether	19.74	20	99%	OK
Hexachlorobenzene	20.88	20	104%	OK
Pentachlorophenol	33.90	40	85%	OK
Phenanthrene	20.24	20	101%	OK
Anthracene	20.85	20	104%	OK
Carbazole	19.92	20	100%	OK
Di-n-butyl phthalate	21.20	20	106%	OK
Fluoranthene	20.81	20	104%	OK
Pyrene	21.53	20	108%	OK
Butyl benzyl phthalate	20.97	20	105%	OK
Benzo[a]anthracene	21.59	20	108%	OK
Chrysene	21.32	20	107%	OK
Bis (2-ethylhexyl) phthalate	20.51	20	103%	OK
di-n-octyl-phthalate	19.28	20	96%	OK
Benzo[b] fluoranthene	20.17	20	101%	OK
Benzo[k] fluoranthene	22.11	20	111%	OK
Benzo[a] pyrene	21.47	20	107%	OK
Indeno [1,2,3-cd] pyrene	19.63	20	98%	OK
Dibenz [a,h] anthracene	20.55	20	103%	OK
Benzo(g,h,i) perylene	20.85	20	104%	OK
2-Fluorophenol	41.34	40	103%	OK
d5-Phenol	30.33	40	76%	OK
d5-nitrobenzene	41.61	40	104%	OK
2-Fluorobiphenyl	41.07	40	103%	OK
2,4,6-Tribromophenol	40.62	40	102%	OK
d14-p-terphenyl	38.69	40	97%	OK

5.26.06

5.26.06 SV

Quantitation Report (Not Reviewed)

Data File : D:\HPCHEM\1\DATA\052406-D\D47730.D Vial: 9
Acq On : 25 May 2006 7:39 am Operator:
Sample : S6857 ABN CC 20/40 Inst : D Instrum
Misc : Multiplr: 1.00
Integration Params: RTEINT.P
Quant Time: May 26 9:37 2006 Quant Results File: S05246.RES

Quant Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Fri May 26 09:35:43 2006
Response via : Initial Calibration
DataAcq Meth : ABN

5.26.06 5.26.06

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) d4-1,4-dichlorobenzene	5.04	152	491009	40.00	ug/ml	0.00
19) d8-Naphthalene	7.95	136	1822322	40.00	ug/ml	0.00
35) d10-Acenaphthene	12.21	164	917674	40.00	ug/ml	0.00
61) d10-Phenanthrene	15.75	188	1589006	40.00	ug/ml	0.00
71) d12-Chrysene	22.17	240	1367670	40.00	ug/ml	0.00
79) d12-Perylene	25.80	264	1331045	40.00	ug/ml	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	2.72	112	780276	41.34	ug/ml	0.00
Spiked Amount 100.000	Range 28 - 110		Recovery = 41.34%			
5) d5-Phenol	4.59	99	686724	30.33	ug/ml	-0.01
Spiked Amount 100.000	Range 22 - 110		Recovery = 30.33%			
20) d5-nitrobenzene	6.36	82	831464	41.61	ug/ml	0.00
Spiked Amount 50.000	Range 45 - 107		Recovery = 83.22%			
39) 2-Fluorobiphenyl	10.65	172	1264418	41.07	ug/ml	0.00
Spiked Amount 50.000	Range 44 - 102		Recovery = 82.14%			
60) 2,4,6-Tribromophenol	14.18	330	199182	40.62	ug/ml	0.00
Spiked Amount 100.000	Range 63 - 127		Recovery = 40.62%#			
73) d14-p-terphenyl	19.70	244	1131010	38.69	ug/ml	0.00
Spiked Amount 50.000	Range 33 - 132		Recovery = 77.38%			
Target Compounds						Qvalue
2) Pyridine	1.02	79	377670	19.60	ug/ml#	100
3) N-nitrosodimethylamine	1.00	42	298998	19.80	ug/ml	98
6) Aniline	4.48	93	97	0.00	ug/ml#	1
7) Phenol	4.62	94	991543	40.30	ug/ml	99
8) 2-Chlorophenol	4.73	128	772745	40.96	ug/ml	98
9) bis(2-chloroethyl) ether	4.59	93	414213	20.88	ug/ml	98
10) 1,3-Dichlorobenzene	4.95	146	390725	21.37	ug/ml	99
11) 1,4-Dichlorobenzene	5.07	146	392520	20.87	ug/ml	97
12) 1,2-Dichlorobenzene	5.47	146	366470	20.71	ug/ml	98
13) Benzyl Alcohol	5.47	108	3627	0.28	ug/ml#	12
14) 2-Methylphenol	5.87	108	714938	40.78	ug/ml	98
15) bis(2-chloroisopropyl)ethe	5.77	45	613661	20.40	ug/ml	98
16) 3+4-Methylphenol	6.23	108	768171	42.15	ug/ml	100
17) Hexachloroethane	6.13	117	184722	20.97	ug/ml	100
18) N-nitroso-di-n-propylamine	6.11	70	262087	22.09	ug/ml	98
21) Nitrobenzene	6.40	123	216753	21.18	ug/ml	98
22) Isophorone	6.94	82	785624	23.56	ug/ml	99
23) 2-Nitrophenol	7.15	139	439882	42.99	ug/ml	93
24) 2,4-Dimethylphenol	7.40	107	750506	40.00	ug/ml	100
25) Bis(2-chloroethoxy)methane	7.54	93	456258	20.48	ug/ml	99

(#) = qualifier out of range (m) = manual integration

Quantitation Report (Not Reviewed)

Data File : D:\HPCHEM\1\DATA\052406-D\D47730.D Vial: 9
Acq On : 25 May 2006 7:39 am Operator:
Sample : S6857 ABN CC 20/40 Inst : D Instrum
Misc : Multiplr: 1.00
Integration Params: RTEINT.P
Quant Time: May 26 9:37 2006 Quant Results File: S05246.RES

Quant Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Fri May 26 09:35:43 2006
Response via : Initial Calibration
DataAcq Meth : ABN

	Compound	R.T.	QIon	Response	Conc Unit	Qvalue
26)	2,4-Dichlorophenol	7.81	162	532352	41.25 ug/ml	98
27)	Benzoic Acid	7.93	105	463640	39.94 ug/ml	99
28)	1,2,4-Trichlorobenzene	7.87	180	287772	20.90 ug/ml	99
29)	Naphthalene	8.00	128	1014125	20.45 ug/ml	99
30)	2,6-Dichlorophenol	8.28	162	535039	40.34 ug/ml	87
31)	4-Chloroaniline	8.28	127	10165	0.46 ug/ml#	1
32)	Hexachlorobutadiene	8.45	225	161915	21.26 ug/ml	99
33)	4-Chloro-3-methylphenol	9.65	107	640428	40.73 ug/ml	99
34)	2-Methylnaphthalene	9.65	141	24270	0.84 ug/ml#	1
36)	Hexachlorocyclopentadiene	10.21	237	77678	25.14 ug/ml	100
37)	2,4,6-Trichlorophenol	10.51	196	352096	41.37 ug/ml	98
38)	2,4,5-Trichlorophenol	10.70	196	349123	38.77 ug/ml	97
40)	2-Chloronaphthalene	10.82	162	556791	20.37 ug/ml	96
41)	2-Nitroaniline	11.25	138	51	0.00 ug/ml#	1
42)	2,6-Dinitrotoluene	11.96	165	167172	21.09 ug/ml	97
43)	Dimethyl Phthalate	11.79	163	657593	20.53 ug/ml	100
44)	Acenaphthylene	11.83	152	908392	21.29 ug/ml	97
45)	3-Nitroaniline	12.28	138	867	0.10 ug/ml#	1
46)	Acenaphthene	12.28	154	504470	20.81 ug/ml	99
47)	2,4-dinitrophenol	12.55	184	182023	39.89 ug/ml	99
48)	Dibenzofuran	12.67	168	1011	0.03 ug/ml#	37
49)	4-Nitrophenol	13.09	109	144162	38.15 ug/ml	96
50)	2,4-Dinitrotoluene	12.89	165	231651	22.46 ug/ml	99
51)	2,3,4,6-Tetrachlorophenol	13.18	232	255825	42.19 ug/ml	98
52)	Fluorene	13.50	166	617979	20.78 ug/ml	99
53)	Diethyl Phthalate	13.52	149	686626	21.68 ug/ml	100
54)	4-Chlorophenyl phenyl eth	13.56	204	290222	20.73 ug/ml#	88
55)	4-Nitroaniline	13.82	138	62	0.01 ug/ml#	1
56)	4,6-Dinitro-2-methylphenol	13.88	198	280405	39.80 ug/ml	97
57)	n-nitrosodiphenylamine	13.91	169	404015	20.09 ug/ml	98
58)	Azobenzene	13.94	77	832223	19.12 ug/ml	96
59)	Biphenyl	10.83	154	1901	0.05 ug/ml#	78
62)	4-bromophenyl phenyl ether	14.73	248	169560	19.74 ug/ml	99
63)	Hexachlorobenzene	15.03	284	207751	20.88 ug/ml	99
64)	Pentachlorophenol	15.56	266	158904	33.90 ug/ml	95
65)	Phenanthrene	15.80	178	810901	20.24 ug/ml	100
66)	Anthracene	15.91	178	842953	20.85 ug/ml	97
67)	Carbazole	16.40	167	778734	19.92 ug/ml	100
68)	Di-n-butyl phthalate	17.45	149	1151809	21.20 ug/ml	99
69)	Fluoranthene	18.68	202	858979	20.81 ug/ml	99
70)	Benzidine	19.13	184	2042	0.10 ug/ml#	93

(#) = qualifier out of range (m) = manual integration

Data File : D:\HPCHEM\1\DATA\052406-D\D47730.D
Acq On : 25 May 2006 7:39 am
Sample : S6857 ABN CC 20/40
Misc :
Integration Params: RTEINT.P
Quant Time: May 26 9:37 2006

Vial: 9
Operator:
Inst : D Instrum
Multiplr: 1.00

Quant Results File: S05246.RES

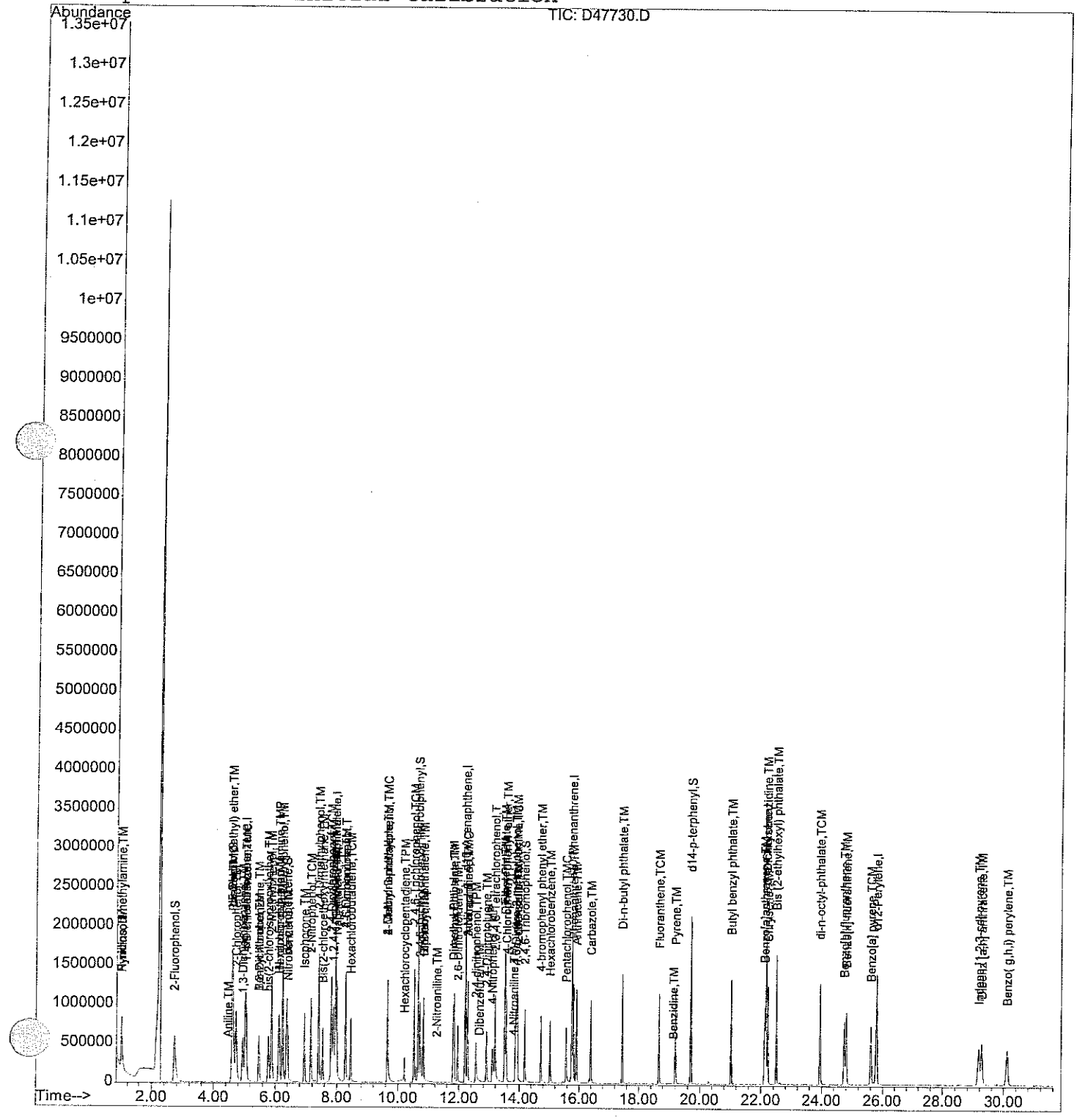
Quant Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Fri May 26 09:35:43 2006
Response via : Initial Calibration
DataAcq Meth : ABN

Compound	R.T.	QIon	Response	Conc Unit	Qvalue
72) Pyrene	19.20	202	886255	21.53 ug/ml	99
74) Butyl benzyl phthalate	21.03	149	514010	20.97 ug/ml	99
75) Benzo[a]anthracene	22.12	228	790647	21.59 ug/ml	99
76) Chrysene	22.22	228	737230	21.32 ug/ml	99
77) 3,3'-Dichlorobenzidine	22.18	252	1045	0.07 ug/ml#	43
78) Bis (2-ethylhexyl) phthala	22.50	149	701181	20.51 ug/ml	98
80) di-n-octyl-phthalate	23.94	149	1131582	19.28 ug/ml	100
81) Benzo[b] fluoranthene	24.74	252	740554	20.17 ug/ml	100
82) Benzo[k] fluoranthene	24.81	252	781929	22.11 ug/ml	98
83) Benzo[a] pyrene	25.62	252	744162	21.47 ug/ml#	91
84) Indeno [1,2,3-cd] pyrene	29.18	276	668136	19.63 ug/ml#	65
85) Dibenz [a,h] anthracene	29.27	278	682101	20.55 ug/ml	99
86) Benzo(g,h,i) perylene	30.12	276	695296	20.85 ug/ml#	85

(#) = qualifier out of range (m) = manual integration

Data File : D:\HPCHEM\1\DATA\052406-D\D47730.D Vial: 9
Acq On : 25 May 2006 7:39 am Operator:
Sample : S6857 ABN CC 20/40 Inst : D Instrum
Misc : Multiplr: 1.00
MS Integration Params: RTEINT.P
Quant Time: May 26 9:37 2006 Quant Results File: S05246.RES

Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Fri May 26 09:35:43 2006
Response via : Initial Calibration



BENZIDINES + ANILINES
CC

Data File Name D47731.D
Date Acquired 5/25/2006 8:12
Sample Name S6858 BENZIDINES + ANILINES

5.26.06

Name	ug/ml Amount	ug/ml Expected	% RECOVERY	70%-130%
Aniline	21.43	20	107%	OK
Benzyl Alcohol	22.63	20	113%	OK
4-Chloroaniline	20.66	20	103%	OK
2-Methylnaphthalene	19.32	20	97%	OK
2-Nitroaniline	18.00	20	90%	OK
3-Nitroaniline	19.04	20	95%	OK
Dibenzofuran	19.26	20	96%	OK
4-Nitroaniline	18.22	20	91%	OK
Benzidine	34.92	20	175%	*
3,3'-Dichlorobenzidine	22.91	20	115%	OK

5.26.06

Quantitation Report (Not Reviewed)

Data File : D:\HPCHEM\1\DATA\052406-D\D47731.D
Acq On : 25 May 2006 8:12 am
Sample : S6858 BENZIDINES + ANILINES CC 20
Misc :
Integration Params: RTEINT.P
Quant Time: May 26 8:45 2006

Vial: 10
Operator:
Inst : D Instrum
Multiplr: 1.00

Quant Results File: S05246.RES

Quant Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Fri May 26 08:10:01 2006
Response via : Initial Calibration
DataAcq Meth : ABN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) d4-1,4-dichlorobenzene	5.03	152	481737	40.00	ug/ml	0.00
19) d8-Naphthalene	7.95	136	1820104	40.00	ug/ml	0.00
35) d10-Acenaphthene	12.21	164	905390	40.00	ug/ml	0.00
61) d10-Phenanthrene	15.74	188	1525849	40.00	ug/ml	0.00
71) d12-Chrysene	22.16	240	1385819	40.00	ug/ml	-0.01
79) d12-Perylene	25.80	264	1359275	40.00	ug/ml	-0.01
System Monitoring Compounds						
4) 2-Fluorophenol	2.73	112	169	0.01	ug/ml	0.00
Spiked Amount 100.000	Range 28 - 110		Recovery = 0.01%#			
5) d5-Phenol	4.61	99	47	0.00	ug/ml	0.00
Spiked Amount 100.000	Range 22 - 110		Recovery = 0.00%#			
20) d5-nitrobenzene	6.36	82	291	0.01	ug/ml	-0.01
Spiked Amount 50.000	Range 45 - 107		Recovery = 0.02%#			
39) 2-Fluorobiphenyl	10.66	172	47	0.00	ug/ml	0.02
Spiked Amount 50.000	Range 44 - 102		Recovery = 0.00%#			
60) 2,4,6-Tribromophenol	14.03	330	47	0.01	ug/ml	-0.15
Spiked Amount 100.000	Range 63 - 127		Recovery = 0.01%#			
73) d14-p-terphenyl	19.64	244	70	0.00	ug/ml	-0.07
Spiked Amount 50.000	Range 33 - 132		Recovery = 0.00%#			
Target Compounds						Qvalue
2) Pyridine	1.02	79	132	0.01	ug/ml#	100
3) N-nitrosodimethylamine	1.02	42	957	0.06	ug/ml#	71
6) Aniline	4.46	93	607160	21.43	ug/ml	99
7) Phenol	4.65	94	452	0.02	ug/ml#	1
8) 2-Chlorophenol	4.74	128	88	0.00	ug/ml#	34
9) bis(2-chloroethyl) ether	4.70	93	943	0.05	ug/ml#	43
10) 1,3-Dichlorobenzene	5.03	146	250	0.01	ug/ml#	1
11) 1,4-Dichlorobenzene	5.03	146	250	0.01	ug/ml#	1
12) 1,2-Dichlorobenzene	5.46	146	71	0.00	ug/ml#	1
13) Benzyl Alcohol	5.46	108	244755	22.63	ug/ml	93
14) 2-Methylphenol	5.88	108	784	0.05	ug/ml#	1
15) bis(2-chloroisopropyl)ethe	5.79	45	52	0.00	ug/ml#	54
16) 3+4-Methylphenol	6.24	108	505	0.03	ug/ml#	76
17) Hexachloroethane	6.12	117	142	0.02	ug/ml#	1
18) N-nitroso-di-n-propylamine	6.12	70	44	0.00	ug/ml#	1
21) Nitrobenzene	6.45	123	44	0.00	ug/ml#	1
22) Isophorone	6.94	82	41	0.00	ug/ml#	1
23) 2-Nitrophenol	7.14	139	110	0.01	ug/ml#	1
24) 2,4-Dimethylphenol	7.38	107	308	0.02	ug/ml#	1
25) Bis(2-chloroethoxy)methane	7.54	93	105	0.00	ug/ml#	20

(#) = qualifier out of range (m) = manual integration

Quantitation Report (Not Reviewed)

Data File : D:\HPCHEM\1\DATA\052406-D\D47731.D Vial: 10
Acq On : 25 May 2006 8:12 am Operator:
Sample : S6858 BENZIDINES + ANILINES CC 20 Inst : D Instrum
Misc : Multiplr: 1.00
Integration Params: RTEINT.P
Quant Time: May 26 8:45 2006 Quant Results File: S05246.RES

Quant Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Fri May 26 08:10:01 2006
Response via : Initial Calibration
DataAcq Meth : ABN

Compound	R.T.	QIon	Response	Conc Unit	Qvalue
26) 2,4-Dichlorophenol	7.83	162	50	0.00 ug/ml#	1
27) Benzoic Acid	7.93	105	98	2.84 ug/ml#	1
28) 1,2,4-Trichlorobenzene	7.88	180	43	0.00 ug/ml#	1
29) Naphthalene	7.99	128	254	0.01 ug/ml#	1
30) 2,6-Dichlorophenol	8.27	162	156	0.01 ug/ml#	1
31) 4-Chloroaniline	8.26	127	448457	20.66 ug/ml	96
33) 4-Chloro-3-methylphenol	9.67	107	40	0.00 ug/ml#	1
34) 2-Methylnaphthalene	9.65	141	534865	19.32 ug/ml#	61
36) Hexachlorocyclopentadiene	10.11	237	54	3.89 ug/ml#	61
37) 2,4,6-Trichlorophenol	10.56	196	44	0.00 ug/ml#	1
38) 2,4,5-Trichlorophenol	10.56	196	44	0.00 ug/ml#	1
40) 2-Chloronaphthalene	10.85	162	50	0.00 ug/ml#	1
41) 2-Nitroaniline	11.26	138	238840	18.00 ug/ml	98
42) 2,6-Dinitrotoluene	11.95	165	64	0.01 ug/ml#	1
43) Dimethyl Phthalate	11.87	163	46	0.00 ug/ml#	1
44) Acenaphthylene	11.83	152	105	0.00 ug/ml#	1
45) 3-Nitroaniline	12.28	138	199230	19.04 ug/ml	95
46) Acenaphthene	12.28	154	76	0.00 ug/ml#	1
47) 2,4-dinitrophenol	12.57	184	100	5.00 ug/ml#	54
48) Dibenzofuran	12.67	168	750746	19.26 ug/ml	99
49) 4-Nitrophenol	13.09	109	168	0.05 ug/ml#	1
50) 2,4-Dinitrotoluene	12.89	165	196	0.02 ug/ml	87
51) 2,3,4,6-Tetrachlorophenol	13.19	232	49	0.01 ug/ml#	44
52) Fluorene	13.51	166	502	0.02 ug/ml#	16
53) Diethyl Phthalate	13.50	149	198	0.01 ug/ml	98
54) 4-Chlorophenyl phenyl eth	13.55	204	55	0.00 ug/ml#	1
55) 4-Nitroaniline	13.78	138	196354	18.22 ug/ml	99
56) 4,6-Dinitro-2-methylphenol	13.87	198	52	3.32 ug/ml#	1
57) n-nitrosodiphenylamine	13.92	169	175	0.01 ug/ml#	24
58) Azobenzene	13.93	77	626	0.01 ug/ml#	1
59) Biphenyl	10.84	154	194	0.01 ug/ml#	34
62) 4-bromophenyl phenyl ether	14.74	248	41	0.00 ug/ml#	1
64) Pentachlorophenol	15.58	266	101	0.02 ug/ml#	74
65) Phenanthrene	15.80	178	461	0.01 ug/ml#	67
66) Anthracene	15.91	178	368	0.01 ug/ml#	77
67) Carbazole	16.39	167	332	0.01 ug/ml#	1
68) Di-n-butyl phthalate	17.44	149	1263	0.02 ug/ml#	78
69) Fluoranthene	18.67	202	82	0.00 ug/ml#	1
70) Benzidine	19.13	184	638587	34.92 ug/ml	99
72) Pyrene	19.21	202	180	0.00 ug/ml#	51
74) Butyl benzyl phthalate	21.02	149	515	0.02 ug/ml	86

(#) = qualifier out of range (m) = manual integration

Quantitation Report (Not Reviewed)

Data File : D:\HPCHEM\1\DATA\052406-D\D47731.D Vial: 10
Acq On : 25 May 2006 8:12 am Operator:
Sample : S6858 BENZIDINES + ANILINES CC 20 Inst : D Instrum
Misc : Multiplr: 1.00
Integration Params: RTEINT.P
Quant Time: May 26 8:45 2006 Quant Results File: S05246.RES

Quant Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Fri May 26 08:10:01 2006
Response via : Initial Calibration
DataAcq Meth : ABN

Compound	R.T.	QIon	Response	Conc Unit	Qvalue
75) Benzo[a]anthracene	22.11	228	100	0.00 ug/ml#	53
76) Chrysene	22.23	228	134	0.00 ug/ml#	51
77) 3,3'-Dichlorobenzidine	22.19	252	333547	22.91 ug/ml	99
78) Bis (2-ethylhexyl) phthala	22.47	149	1122	0.03 ug/ml#	87
80) di-n-octyl-phthalate	23.96	149	263	0.00 ug/ml#	1
81) Benzo[b] fluoranthene	24.74	252	416	0.01 ug/ml#	85
82) Benzo[k] fluoranthene	24.80	252	206	0.01 ug/ml#	1
83) Benzo[a] pyrene	25.62	252	83	0.00 ug/ml#	54
84) Indeno [1,2,3-cd] pyrene	29.18	276	52	0.00 ug/ml#	62
85) Dibenz [a,h] anthracene	29.24	278	248	0.01 ug/ml#	52
86) Benzo(g,h,i) perylene	30.15	276	91	0.00 ug/ml#	78

(#) = qualifier out of range (m) = manual integration

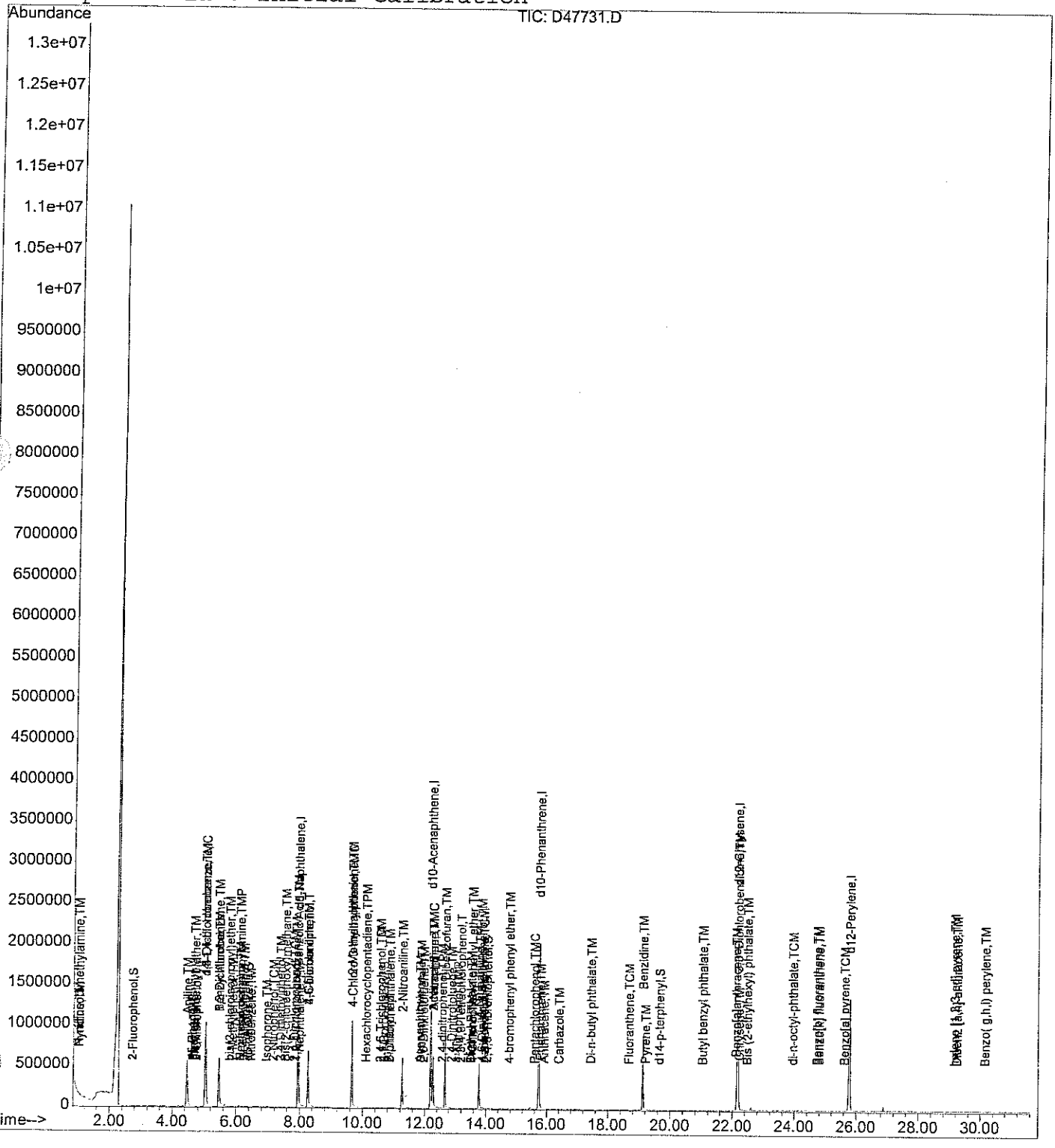
Quantitation Report

Data File : D:\HPCHEM\1\DATA\052406-D\D47731.D
Acq On : 25 May 2006 8:12 am
Sample : S6858 BENZIDINES + ANILINES CC 20
Misc :
MS Integration Params: RTEINT.P
Quant Time: May 26 8:45 2006

Vial: 10
Operator:
Inst : D Instrum
Multiplr: 1.00

Quant Results File: S05246.RES

Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Fri May 26 09:35:43 2006
Response via : Initial Calibration



Evaluate Continuing Calibration Report

Data File : D:\HPCHEM\1\DATA\053006-D\D47756SC.D Vial: 2
Acq On : 30 May 2006 7:24 am Operator:
Sample : S6849 PT3 20/40 Inst : D Instrum
Misc : Multiplr: 1.00
MS Integration Params: RTEINT.P

Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Tue May 30 07:32:16 2006
Response via : Multiple Level Calibration

Min. RRF : 0.050 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
Max. RRF Dev : 20% Max. Rel. Area : 200%

		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
1	I	d4-1,4-dichlorobenzene	1.000	1.000	0.0	119	0.00
2	TM	Pyridine	1.570	1.520	3.2	114	0.00
3	TM	N-nitrosodimethylamine	1.230	1.223	0.6	115	0.00
4	S	2-Fluorophenol	1.538	1.535	0.2	117	0.00
5	S	d5-Phenol	1.844	1.831	0.7	117	0.00
6	TM	Aniline	2.438	2.172	10.9	104	0.00
7	TMC	Phenol	2.004	1.956	2.4	117	0.00
8	TM	2-Chlorophenol	1.537	1.525	0.8	118	0.00
9	TM	bis(2-chloroethyl) ether	1.616	1.604	0.7	118	0.00
10	TM	1,3-Dichlorobenzene	1.489	1.492	-0.2	119	0.00
11	TMC	1,4-Dichlorobenzene	1.532	1.541	0.6	116	0.00
12	TM	1,2-Dichlorobenzene	1.441	1.449	-0.6	118	0.00
13	TM	Benzyl Alcohol	1.056	1.060	-0.4	114	0.00
14	TM	2-Methylphenol	1.428	1.413	1.1	116	0.00
15	TM	bis(2-chloroisopropyl)ether	2.451	2.270	7.4	108	0.00
16	TM	3+4-Methylphenol	1.485	1.513	-1.9	118	0.00
17	TM	Hexachloroethane	0.718	0.737	-2.6	122	0.00
18	TMP	N-nitroso-di-n-propylamine	0.966	0.919	4.9	114	0.00
19	I	d8-Naphthalene	1.000	1.000	0.0	119	0.00
20	S	d5-nitrobenzene	0.439	0.455	-3.6	118	0.00
21	TM	Nitrobenzene	0.225	0.227	-0.9	116	0.00
22	TM	Isophorone	0.732	0.736	-0.5	116	0.00
23	TCM	2-Nitrophenol	0.225	0.233	-3.6	122	0.00
24	TM	2,4-Dimethylphenol	0.412	0.393	4.6	111	0.00
25	TM	Bis(2-chloroethoxy)methane	0.489	0.490	-0.2	117	0.00
26	TCM	2,4-Dichlorophenol	0.283	0.297	-4.9	122	0.00
27	TM	Benzoic Acid	0.215	0.203	5.6	103	0.00
28	TM	1,2,4-Trichlorobenzene	0.302	0.319	-5.6	125	0.00
29	TM	Naphthalene	1.088	1.111	-2.1	120	0.00
30	T	2,6-Dichlorophenol	0.291	0.294	-1.0	121	0.00
31	TM	4-Chloroaniline	0.488	0.487	0.2	119	0.00
32	TCM	Hexachlorobutadiene	0.167	0.177	-6.0	125	0.00
33	TMC	4-Chloro-3-methylphenol	0.345	0.368	-6.7	124	0.00
34	TM	2-Methylnaphthalene	0.633	0.641	-1.3	118	0.00
35	I	d10-Acenaphthene	1.000	1.000	0.0	124	0.00
36	TPM	Hexachlorocyclopentadiene	0.122	0.115	5.7	132	0.00
37	TCM	2,4,6-Trichlorophenol	0.371	0.375	-1.1	121	0.00
38	TM	2,4,5-Trichlorophenol	0.393	0.391	0.5	126	0.00
39	S	2-Fluorobiphenyl	1.342	1.330	0.9	123	0.00

(#) = Out of Range

Evaluate Continuing Calibration Report

Data File : D:\HPCHEM\1\DATA\053006-D\D47756SC.D Vial: 2
Acq On : 30 May 2006 7:24 am Operator:
Sample : S6849 PT3 20/40 Inst : D Instrum
Misc : Multiplr: 1.00
MS Integration Params: RTEINT.P

Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Tue May 30 07:32:16 2006
Response via : Multiple Level Calibration

Min. RRF : 0.050 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
Max. RRF Dev : 20% Max. Rel. Area : 200%

		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
40	TM	2-Chloronaphthalene	1.191	1.166	2.1	123	0.00
41	TM	2-Nitroaniline	0.502	0.517	-3.0	121	0.00
42	TM	2,6-Dinitrotoluene	0.345	0.359	-4.1	125	0.00
43	TM	Dimethyl Phthalate	1.396	1.400	-0.3	124	0.00
44	TM	Acenaphthylene	1.860	1.838	1.2	120	0.00
45	TM	3-Nitroaniline	0.397	0.414	-4.3	125	0.00
46	TMC	Acenaphthene	1.057	1.062	-0.5	125	0.00
47	TPM	2,4-dinitrophenol	0.167	0.184	-10.2	121	0.00
48	TM	Dibenzofuran	1.661	1.688	-1.6	127	0.00
49	TMP	4-Nitrophenol	0.165	0.160	3.0	128	0.00
50	TM	2,4-Dinitrotoluene	0.450	0.480	-6.7	126	0.00
51	T	2,3,4,6-Tetrachlorophenol	0.250	0.259	-3.6	124	0.00
52	TM	Fluorene	1.296	1.302	-0.5	124	0.00
53	TM	Diethyl Phthalate	1.380	1.408	-2.0	123	0.00
54	TM	4-Chlorophenyl phenyl ethe	0.610	0.637	-4.4	129	0.00
55	TM	4-Nitroaniline	0.403	0.408	-1.2	117	0.00
56	TM	4,6-Dinitro-2-methylphenol	0.266	0.301	-13.2	126	0.00
57	TCM	n-nitrosodiphenylamine	0.876	0.884	-0.9	125	0.00
58	TM	Azobenzene	1.897	1.870	1.4	120	0.00
59	TM	Biphenyl	1.513	1.486	1.8	124	0.00
60	S	2,4,6-Tribromophenol	0.214	0.223	-4.2	129	0.00
61	I	d10-Phenanthrene	1.000	1.000	0.0	130	0.00
62	TM	4-bromophenyl phenyl ether	0.216	0.214	0.9	130	0.00
63	TM	Hexachlorobenzene	0.251	0.247	1.6	124	0.00
64	TMC	Pentachlorophenol	0.117	0.106	9.4	124	0.00
65	TM	Phenanthrene	1.009	0.985	2.4	128	0.00
66	TM	Anthracene	1.018	0.990	2.8	122	0.00
67	TM	Carbazole	0.984	0.946	3.9	123	0.00
68	TM	Di-n-butyl phthalate	1.367	1.368	-0.1	123	0.00
69	TCM	Fluoranthene	1.039	1.066	-2.6	130	0.00
70	TM	Benzidine	0.507	0.154	69.6#	37#	0.00
71	I	d12-Chrysene	1.000	1.000	0.0	132	0.00
72	TM	Pyrene	1.204	1.220	-1.3	132	0.00
73	S	d14-p-terphenyl	0.855	0.838	2.0	130	0.00
74	TM	Butyl benzyl phthalate	0.717	0.704	1.8	126	0.00
75	TM	Benzo[a]anthracene	1.071	1.084	-1.2	132	0.00
76	TM	Chrysene	1.011	1.004	0.7	129	0.00
77	TM	3,3'-Dichlorobenzidine	0.455	0.480	-5.5	138	0.00
78	TM	Bis (2-ethylhexyl) phthalat	1.000	0.971	2.9	129	0.00

(#) = Out of Range

Evaluate Continuing Calibration Report

Data File : D:\HPCHEM\1\DATA\053006-D\D47756SC.D Vial: 2
Acq On : 30 May 2006 7:24 am Operator:
Sample : S6849 PT3 20/40 Inst : D Instrum
Misc : Multiplr: 1.00
MS Integration Params: RTEINT.P

Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Tue May 30 07:32:16 2006
Response via : Multiple Level Calibration

Min. RRF : 0.050 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
Max. RRF Dev : 20% Max. Rel. Area : 200%

		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
79	I	d12-Perylene	1.000	1.000	0.0	145	0.00
80	TCM	di-n-octyl-phthalate	1.764	1.609	8.8	129	0.00
81	TM	Benzo[b] fluoranthene	1.103	1.042	5.5	136	0.00
82	TM	Benzo[k] fluoranthene	1.063	1.050	1.2	140	0.00
83	TCM	Benzo[a] pyrene	1.042	1.003	3.7	137	0.00
84	TM	Indeno [1,2,3-cd] pyrene	1.023	1.061	-3.7	151	0.00
85	TM	Dibenz [a,h] anthracene	0.998	1.054	-5.6	153	0.00
86	TM	Benzo(g,h,i) perylene	1.002	1.035	-3.3	149	0.00

Quantitation Report (Not Reviewed)

Data File : D:\HPCHEM\1\DATA\053006-D\D47757SC.D Vial: 3
Acq On : 30 May 2006 7:57 am Operator:
Sample : S6760 B+A 20 Inst : D Instrum
Misc : Multiplr: 1.00
Integration Params: RTEINT.P
Quant Time: May 30 8:10 2006 Quant Results File: S05246.RES

Quant Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Tue May 30 07:32:16 2006
Response via : Initial Calibration
DataAcq Meth : ABN

5.31.06 SW

5/30/06

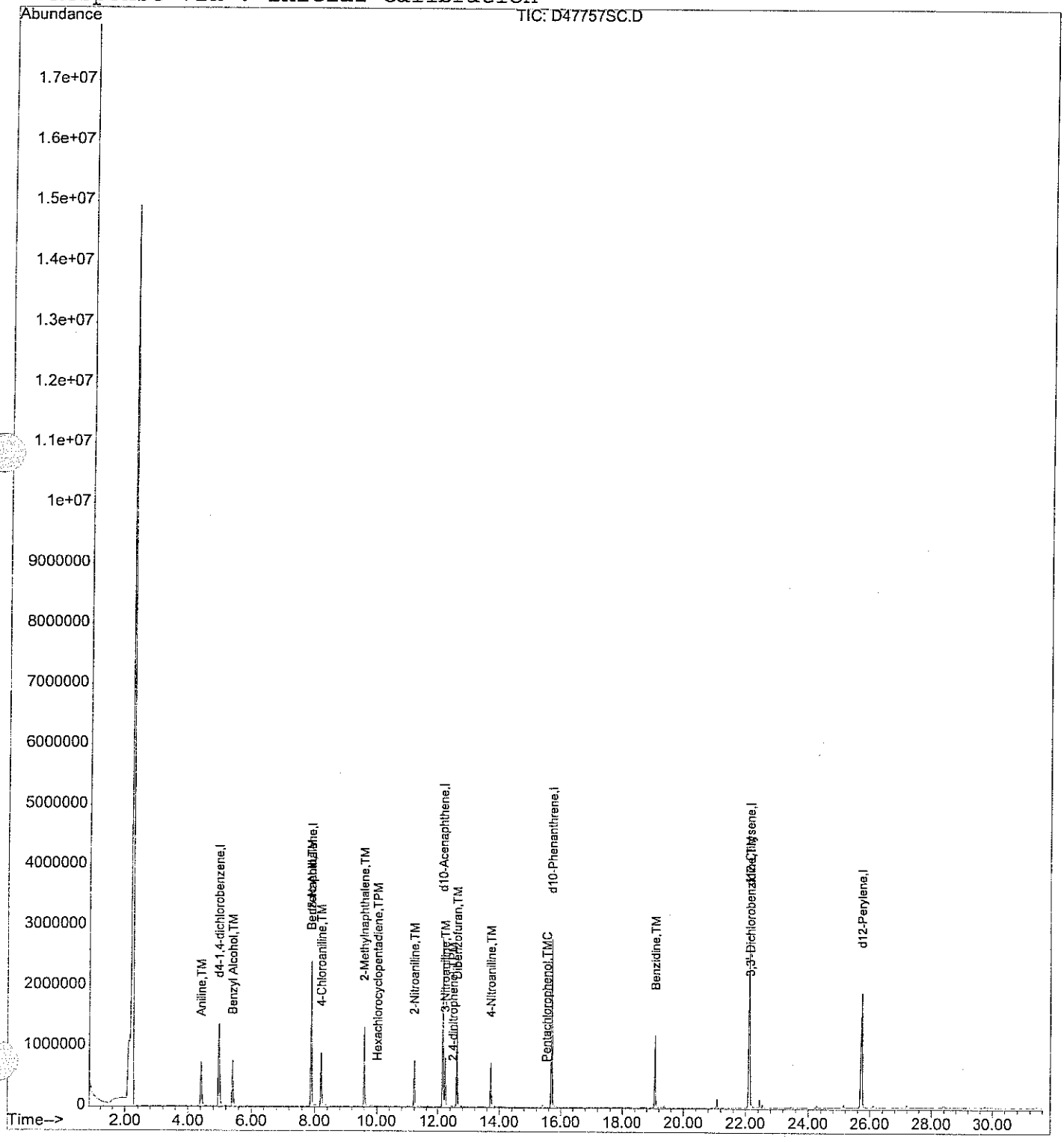
Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) d4-1,4-dichlorobenzene	4.98	152	620189	40.00	ug/ml	0.00
19) d8-Naphthalene	7.89	136	2352118	40.00	ug/ml	0.00
35) d10-Acenaphthene	12.15	164	1202449	40.00	ug/ml	0.00
61) d10-Phenanthrene	15.69	188	2170249	40.00	ug/ml	0.00
71) d12-Chrysene	22.11	240	1951558	40.00	ug/ml	-0.01
79) d12-Perylene	25.73	264	2018812	40.00	ug/ml	-0.01
System Monitoring Compounds						
4) 2-Fluorophenol	2.68	112	199	0.01	ug/ml	0.00
Spiked Amount 100.000	Range 28 - 110		Recovery =	0.01%#		
5) d5-Phenol	4.56	99	44	0.00	ug/ml	0.01
Spiked Amount 100.000	Range 22 - 110		Recovery =	0.00%#		
20) d5-nitrobenzene	6.32	82	103	0.00	ug/ml	0.00
Spiked Amount 50.000	Range 45 - 107		Recovery =	0.00%#		
39) 2-Fluorobiphenyl	10.66	172	261	0.01	ug/ml	0.06
Spiked Amount 50.000	Range 44 - 102		Recovery =	0.02%#		
60) 2,4,6-Tribromophenol	0.00	330	0	0.00	ug/ml	
Spiked Amount 100.000	Range 63 - 127		Recovery =	0.00%#		
73) d14-p-terphenyl	19.67	244	50	0.00	ug/ml	0.01
Spiked Amount 50.000	Range 33 - 132		Recovery =	0.00%#		
Target Compounds						Qvalue
6) Aniline	4.41	93	741588	19.62	ug/ml	98
13) Benzyl Alcohol	5.41	108	310405	18.95	ug/ml	97
27) Benzoic Acid	7.88	105	540	5.21	ug/ml#	1
31) 4-Chloroaniline	8.21	127	567395	19.75	ug/ml	95
34) 2-Methylnaphthalene	9.60	141	651203	17.49	ug/ml#	58
36) Hexachlorocyclopentadiene	10.02	237	53	5.48	ug/ml#	21
41) 2-Nitroaniline	11.20	138	299635	19.84	ug/ml	100
45) 3-Nitroaniline	12.23	138	245881	20.58	ug/ml	96
47) 2,4-dinitrophenol	12.49	184	70	5.39	ug/ml#	60
48) Dibenzofuran	12.61	168	1005387	20.14	ug/ml	100
55) 4-Nitroaniline	13.73	138	252603	20.86	ug/ml	99
64) Pentachlorophenol	15.53	266	65	6.50	ug/ml#	25
70) Benzidine	19.07	184	818901	29.76	ug/ml	98
77) 3,3'-Dichlorobenzidine	22.14	252	426182	19.21	ug/ml	99

(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : D:\HPCHEM\1\DATA\053006-D\D47757SC.D Vial: 3
Acq On : 30 May 2006 7:57 am Operator:
Sample : S6760 B+A 20 Inst : D Instrum
Misc : Multiplr: 1.00
MS Integration Params: RTEINT.P
Quant Time: May 30 8:10 2006 Quant Results File: S05246.RES

Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Tue May 30 07:32:16 2006
Response via : Initial Calibration



Evaluate Continuing Calibration Report

Data File : D:\HPCHEM\1\DATA\053006-D\D47772SC.D Vial: 2
Acq On : 30 May 2006 9:08 pm Operator:
Sample : S6849 PT3 20/40 Inst : D Instrum
Misc : Multiplr: 1.00
MS Integration Params: RTEINT.P

Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Tue May 30 20:31:56 2006
Response via : Multiple Level Calibration

5-31-06 SL

Min. RRF : 0.050 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
Max. RRF Dev : 20% Max. Rel. Area : 200%

		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
1	I	d4-1,4-dichlorobenzene	1.000	1.000	0.0	122	-0.01
2	TM	Pyridine	1.570	1.493	4.9	115	-0.01
3	TM	N-nitrosodimethylamine	1.230	1.225	0.4	118	0.00
4	S	2-Fluorophenol	1.538	1.500	2.5	117	-0.02
5	S	d5-Phenol	1.844	1.808	2.0	118	-0.02
6	TM	Aniline	2.438	2.088	14.4	102	-0.01
7	TMC	Phenol	2.004	1.933	3.5	118	-0.01
8	TM	2-Chlorophenol	1.537	1.524	0.8	121	-0.01
9	TM	bis(2-chloroethyl) ether	1.616	1.641	-1.5	124	-0.01
10	TM	1,3-Dichlorobenzene	1.489	1.486	0.2	122	-0.01
11	TMC	1,4-Dichlorobenzene	1.532	1.595	-4.1	123	-0.01
12	TM	1,2-Dichlorobenzene	1.441	1.465	-1.7	122	0.00
13	TM	Benzyl Alcohol	1.056	1.051	0.5	115	-0.01
14	TM	2-Methylphenol	1.428	1.390	2.7	117	-0.02
15	TM	bis(2-chloroisopropyl)ether	2.451	2.275	7.2	111	0.00
16	TM	3+4-Methylphenol	1.485	1.465	1.3	117	-0.01
17	TM	Hexachloroethane	0.718	0.735	-2.4	124	-0.01
18	TMP	N-nitroso-di-n-propylamine	0.966	0.922	4.6	117	0.00
19	I	d8-Naphthalene	1.000	1.000	0.0	121	0.00
20	S	d5-nitrobenzene	0.439	0.449	-2.3	119	0.00
21	TM	Nitrobenzene	0.225	0.226	-0.4	118	0.00
22	TM	Isophorone	0.732	0.713	2.6	115	-0.01
23	TCM	2-Nitrophenol	0.225	0.230	-2.2	122	-0.01
24	TM	2,4-Dimethylphenol	0.412	0.394	4.4	113	-0.01
25	TM	Bis(2-chloroethoxy)methane	0.489	0.485	0.8	118	0.00
26	TCM	2,4-Dichlorophenol	0.283	0.293	-3.5	123	-0.01
27	TM	Benzoic Acid	0.215	0.216	-0.5	112	-0.02
28	TM	1,2,4-Trichlorobenzene	0.302	0.321	-6.3	129	0.00
29	TM	Naphthalene	1.088	1.094	-0.6	120	-0.01
30	T	2,6-Dichlorophenol	0.291	0.287	1.4	120	-0.01
31	TM	4-Chloroaniline	0.488	0.468	4.1	116	0.00
32	TCM	Hexachlorobutadiene	0.167	0.183	-9.6	132	-0.01
33	TMC	4-Chloro-3-methylphenol	0.345	0.354	-2.6	122	-0.01
34	TM	2-Methylnaphthalene	0.633	0.630	0.5	118	-0.01
35	I	d10-Acenaphthene	1.000	1.000	0.0	124	0.00
36	TPM	Hexachlorocyclopentadiene	0.122	0.104	14.8	119	0.00
37	TCM	2,4,6-Trichlorophenol	0.371	0.380	-2.4	123	-0.01
38	TM	2,4,5-Trichlorophenol	0.393	0.378	3.8	121	-0.02
39	S	2-Fluorobiphenyl	1.342	1.321	1.6	122	-0.01

(#) = Out of Range

Evaluate Continuing Calibration Report

Data File : D:\HPCHEM\1\DATA\053006-D\D47772SC.D Vial: 2
Acq On : 30 May 2006 9:08 pm Operator:
Sample : S6849 PT3 20/40 Inst : D Instrum
Misc : Multiplr: 1.00
MS Integration Params: RTEINT.P

Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Tue May 30 20:31:56 2006
Response via : Multiple Level Calibration

Min. RRF : 0.050 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
Max. RRF Dev : 20% Max. Rel. Area : 200%

		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
40	TM	2-Chloronaphthalene	1.191	1.165	2.2	123	0.00
41	TM	2-Nitroaniline	0.502	0.515	-2.6	120	0.00
42	TM	2,6-Dinitrotoluene	0.345	0.357	-3.5	124	0.00
43	TM	Dimethyl Phthalate	1.396	1.393	0.2	123	-0.01
44	TM	Acenaphthylene	1.860	1.819	2.2	118	-0.01
45	TM	3-Nitroaniline	0.397	0.405	-2.0	122	0.00
46	TMC	Acenaphthene	1.057	1.051	0.6	123	-0.01
47	TPM	2,4-dinitrophenol	0.167	0.183	-9.6	120	-0.01
48	TM	Dibenzofuran	1.661	1.681	-1.2	126	0.00
49	TMP	4-Nitrophenol	0.165	0.160	3.0	127	-0.02
50	TM	2,4-Dinitrotoluene	0.450	0.460	-2.2	120	0.00
51	T	2,3,4,6-Tetrachlorophenol	0.250	0.260	-4.0	124	-0.01
52	TM	Fluorene	1.296	1.300	-0.3	123	0.00
[illegible]	TM	Diethyl Phthalate	1.380	1.412	-2.3	123	-0.01
54	TM	4-Chlorophenyl phenyl ethe	0.610	0.630	-3.3	127	0.00
55	TM	4-Nitroaniline	0.403	0.405	-0.5	116	0.00
56	TM	4,6-Dinitro-2-methylphenol	0.266	0.287	-7.9	119	-0.01
57	TCM	n-nitrosodiphenylamine	0.876	0.876	0.0	123	0.00
58	TM	Azobenzene	1.897	1.846	2.7	118	0.00
59	TM	Biphenyl	1.513	1.476	2.4	123	0.00
60	S	2,4,6-Tribromophenol	0.214	0.205	4.2	118	0.00
61	I	d10-Phenanthrene	1.000	1.000	0.0	126	0.00
62	TM	4-bromophenyl phenyl ether	0.216	0.215	0.5	127	-0.01
63	TM	Hexachlorobenzene	0.251	0.249	0.8	120	-0.01
64	TMC	Pentachlorophenol	0.117	0.116	0.9	131	-0.01
65	TM	Phenanthrene	1.009	0.995	1.4	125	-0.01
66	TM	Anthracene	1.018	1.016	0.2	121	0.00
67	TM	Carbazole	0.984	0.969	1.5	122	0.00
68	TM	Di-n-butyl phthalate	1.367	1.391	-1.8	122	-0.01
69	TCM	Fluoranthene	1.039	1.071	-3.1	127	-0.01
70	TM	Benzidine	0.507	0.145	71.4#	34#	0.00
71	I	d12-Chrysene	1.000	1.000	0.0	121	-0.01
72	TM	Pyrene	1.204	1.296	-7.6	129	-0.01
73	S	d14-p-terphenyl	0.855	0.872	-2.0	124	-0.01
74	TM	Butyl benzyl phthalate	0.717	0.742	-3.5	122	0.00
75	TM	Benzo[a]anthracene	1.071	1.098	-2.5	123	-0.01
[illegible]	TM	Chrysene	1.011	1.036	-2.5	122	-0.01
[illegible]	TM	3,3'-Dichlorobenzidine	0.455	0.461	-1.3	122	0.00
78	TM	Bis (2-ethylhexyl) phthalat	1.000	1.014	-1.4	123	0.00

(#) = Out of Range

Evaluate Continuing Calibration Report

Data File : D:\HPCHEM\1\DATA\053006-D\D47772SC.D Vial: 2
Acq On : 30 May 2006 9:08 pm Operator:
Sample : S6849 PT3 20/40 Inst : D Instrum
Misc : Multiplr: 1.00
MS Integration Params: RTEINT.P

Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Tue May 30 20:31:56 2006
Response via : Multiple Level Calibration

Min. RRF : 0.050 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
Max. RRF Dev : 20% Max. Rel. Area : 200%

		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
79	I	d12-Perylene	1.000	1.000	0.0	118	-0.03
80	TCM	di-n-octyl-phthalate	1.764	1.863	-5.6	122	-0.01
81	TM	Benzo[b] fluoranthene	1.103	1.126	-2.1	119	-0.01
82	TM	Benzo[k] fluoranthene	1.063	1.103	-3.8	120	-0.02
83	TCM	Benzo[a] pyrene	1.042	1.016	2.5	113	-0.02
84	TM	Indeno [1,2,3-cd] pyrene	1.023	0.954	6.7	110	-0.03
85	TM	Dibenz [a,h] anthracene	0.998	0.934	6.4	110	-0.03
86	TM	Benzo(g,h,i) perylene	1.002	0.928	7.4	108	-0.04

Quantitation Report (Not Reviewed)

Data File : D:\HPCHEM\1\DATA\053006-D\D47773SC.D Vial: 3
Acq On : 30 May 2006 9:46 pm Operator:
Sample : S6760 B+A 20 Inst : D Instrum
Misc : Multiplr: 1.00
Integration Params: RTEINT.P
Quant Time: May 30 22:29 2006 Quant Results File: S05246.RES

Quant Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Tue May 30 20:31:56 2006
Response via : Initial Calibration
DataAcq Meth : ABN

5.31.06 ✓

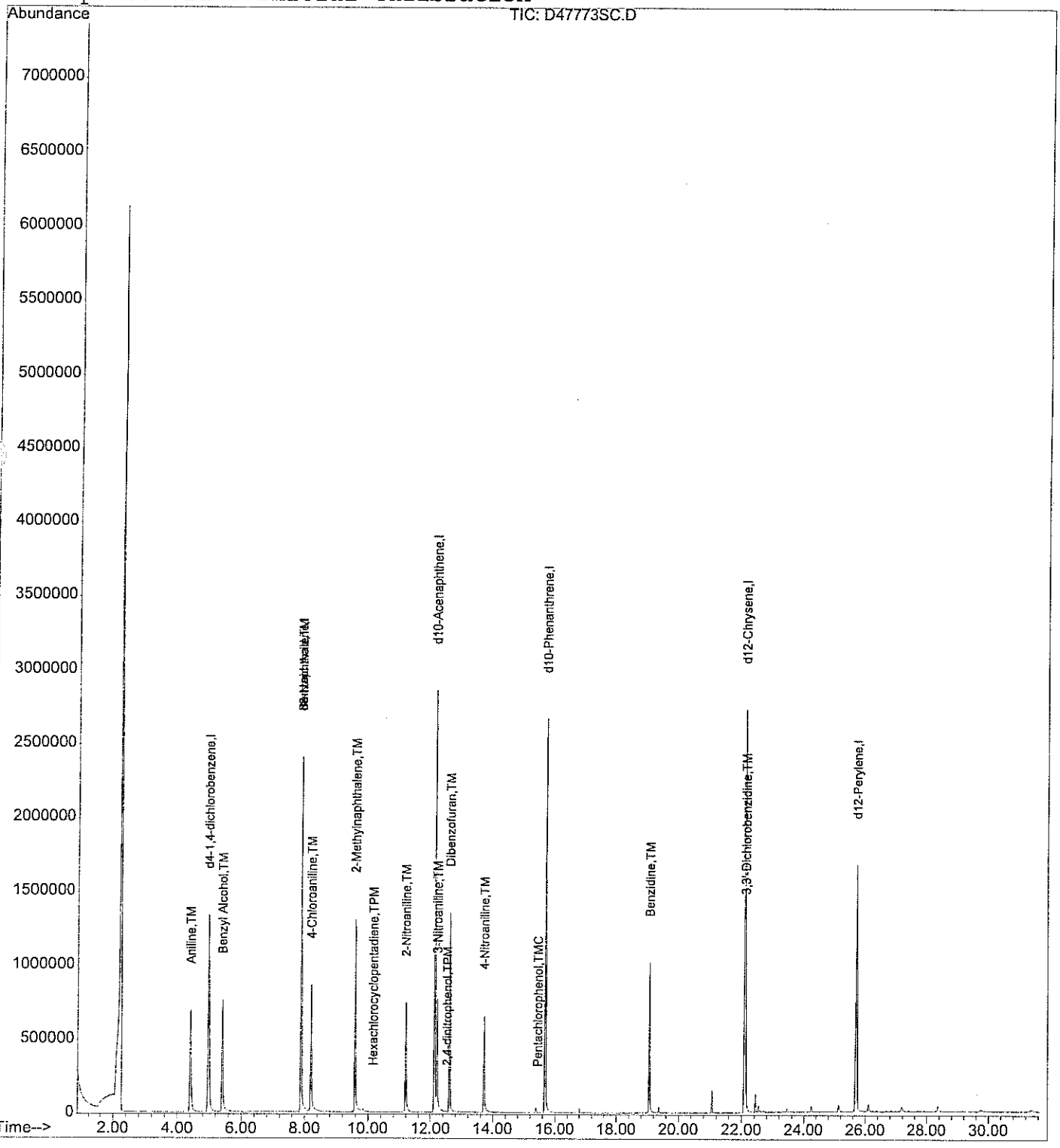
Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) d4-1,4-dichlorobenzene	4.97	152	631919	40.00	ug/ml	-0.01
19) d8-Naphthalene	7.88	136	2372906	40.00	ug/ml	-0.01
35) d10-Acenaphthene	12.14	164	1211648	40.00	ug/ml	-0.01
61) d10-Phenanthrene	15.69	188	2126590	40.00	ug/ml	-0.01
71) d12-Chrysene	22.10	240	1795092	40.00	ug/ml	-0.02
79) d12-Perylene	25.71	264	1664861	40.00	ug/ml	-0.03
System Monitoring Compounds						
4) 2-Fluorophenol	2.70	112	47	0.00	ug/ml	0.01
Spiked Amount 75.000	Range 28 - 110		Recovery =	0.00%#		
5) d5-Phenol	4.55	99	95	0.00	ug/ml	0.00
Spiked Amount 75.000	Range 22 - 110		Recovery =	0.00%#		
20) d5-nitrobenzene	6.29	82	144	0.01	ug/ml	-0.03
Spiked Amount 50.000	Range 45 - 107		Recovery =	0.02%#		
39) 2-Fluorobiphenyl	10.63	172	187	0.00	ug/ml	0.04
Spiked Amount 50.000	Range 44 - 102		Recovery =	0.00%#		
60) 2,4,6-Tribromophenol	0.00	330	0	0.00	ug/ml	
Spiked Amount 75.000	Range 63 - 127		Recovery =	0.00%#		
73) d14-p-terphenyl	19.68	244	51	0.00	ug/ml	0.02
Spiked Amount 50.000	Range 33 - 132		Recovery =	0.00%#		
Target Compounds						Qvalue
6) Aniline	4.41	93	747108	19.40	ug/ml	98
13) Benzyl Alcohol	5.40	108	318256	19.07	ug/ml	99
27) Benzoic Acid	7.88	105	534	5.21	ug/ml#	1
31) 4-Chloroaniline	8.20	127	558103	19.26	ug/ml	94
34) 2-Methylnaphthalene	9.59	141	647813	17.24	ug/ml#	58
36) Hexachlorocyclopentadiene	10.15	237	132	5.49	ug/ml	97
41) 2-Nitroaniline	11.20	138	302193	19.86	ug/ml	97
45) 3-Nitroaniline	12.22	138	239085	19.86	ug/ml	95
47) 2,4-dinitrophenol	12.52	184	194	5.41	ug/ml#	46
48) Dibenzofuran	12.61	168	984937	19.58	ug/ml	99
55) 4-Nitroaniline	13.72	138	237987	19.50	ug/ml	98
64) Pentachlorophenol	15.44	266	57	6.50	ug/ml#	28
70) Benzidine	19.06	184	739023	27.41	ug/ml	99
77) 3,3'-Dichlorobenzidine	22.13	252	376142	18.44	ug/ml	99

(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : D:\HPCHEM\1\DATA\053006-D\D47773SC.D Vial: 3
Acq On : 30 May 2006 9:46 pm Operator:
Sample : S6760 B+A 20 Inst : D Instrum
Misc : Multiplr: 1.00
MS Integration Params: RTEINT.P
Quant Time: May 30 22:29 2006 Quant Results File: S05246.RES

Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Tue May 30 20:31:56 2006
Response via : Initial Calibration



Evaluate Continuing Calibration Report

Data File : D:\HPCHEM\1\DATA\053106-D\D47795SC.D Vial: 2
Acq On : 31 May 2006 11:03 am Operator:
Sample : S6849 PT3 20/40 Inst : D Instrum
Misc : Multiplr: 1.00
MS Integration Params: RTEINT.P

Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Wed May 31 12:09:53 2006
Response via : Multiple Level Calibration

6-1-06 SL ✓ 5/31/06

Min. RRF : 0.050 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
Max. RRF Dev : 20% Max. Rel. Area : 200%

		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
1	I	d4-1,4-dichlorobenzene	1.000	1.000	0.0	112	0.00
2	TM	Pyridine	1.570	1.476	6.0	104	0.00
3	TM	N-nitrosodimethylamine	1.230	1.194	2.9	106	0.00
4	S	2-Fluorophenol	1.538	1.542	-0.3	110	0.00
5	S	d5-Phenol	1.844	1.824	1.1	109	0.00
6	TM	Aniline	2.438	2.128	12.7	95	0.00
7	TMC	Phenol	2.004	1.965	1.9	110	0.00
8	TM	2-Chlorophenol	1.537	1.520	1.1	111	0.00
9	TM	bis(2-chloroethyl) ether	1.616	1.520	5.9	105	0.00
10	TM	1,3-Dichlorobenzene	1.489	1.475	0.9	111	0.00
11	TMC	1,4-Dichlorobenzene	1.532	1.545	-0.8	110	0.00
12	TM	1,2-Dichlorobenzene	1.441	1.408	2.3	107	0.00
13	TM	Benzyl Alcohol	1.056	1.060	-0.4	106	0.00
[illegible]	TM	2-Methylphenol	1.428	1.410	1.3	109	0.00
[illegible]	TM	bis(2-chloroisopropyl)ether	2.451	2.243	8.5	100	0.00
16	TM	3+4-Methylphenol	1.485	1.476	0.6	108	0.00
17	TM	Hexachloroethane	0.718	0.719	-0.1	112	0.00
18	TMP	N-nitroso-di-n-propylamine	0.966	0.904	6.4	105	0.00
19	I	d8-Naphthalene	1.000	1.000	0.0	110	0.00
20	S	d5-nitrobenzene	0.439	0.447	-1.8	108	0.00
21	TM	Nitrobenzene	0.225	0.229	-1.8	108	0.00
22	TM	Isophorone	0.732	0.720	1.6	105	0.00
23	TCM	2-Nitrophenol	0.225	0.237	-5.3	115	0.00
24	TM	2,4-Dimethylphenol	0.412	0.389	5.6	102	0.00
25	TM	Bis(2-chloroethoxy)methane	0.489	0.495	-1.2	109	0.00
26	TCM	2,4-Dichlorophenol	0.283	0.298	-5.3	113	0.00
27	TM	Benzoic Acid	0.215	0.249	-15.8	117	0.00
28	TM	1,2,4-Trichlorobenzene	0.302	0.312	-3.3	114	0.00
29	TM	Naphthalene	1.088	1.102	-1.3	110	0.00
30	T	2,6-Dichlorophenol	0.291	0.293	-0.7	111	0.00
31	TM	4-Chloroaniline	0.488	0.475	2.7	107	0.00
32	TCM	Hexachlorobutadiene	0.167	0.175	-4.8	115	0.00
33	TMC	4-Chloro-3-methylphenol	0.345	0.359	-4.1	112	0.00
34	TM	2-Methylnaphthalene	0.633	0.636	-0.5	108	0.00
35	I	d10-Acenaphthene	1.000	1.000	0.0	109	0.00
36	TPM	Hexachlorocyclopentadiene	0.122	0.104	14.8	105	0.00
[illegible]	TCM	2,4,6-Trichlorophenol	0.371	0.386	-4.0	110	0.00
[illegible]	TM	2,4,5-Trichlorophenol	0.393	0.405	-3.1	115	0.00
39	S	2-Fluorobiphenyl	1.342	1.352	-0.7	110	0.00

(#) = Out of Range

Evaluate Continuing Calibration Report

Data File : D:\HPCHEM\1\DATA\053106-D\D47795SC.D Vial: 2
Acq On : 31 May 2006 11:03 am Operator:
Sample : S6849 PT3 20/40 Inst : D Instrum
Misc : Multiplr: 1.00
MS Integration Params: RTEINT.P

Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Wed May 31 12:09:53 2006
Response via : Multiple Level Calibration

Min. RRF : 0.050 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
Max. RRF Dev : 20% Max. Rel. Area : 200%

		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
40	TM	2-Chloronaphthalene	1.191	1.179	1.0	110	0.00
41	TM	2-Nitroaniline	0.502	0.528	-5.2	109	0.00
42	TM	2,6-Dinitrotoluene	0.345	0.356	-3.2	110	0.00
43	TM	Dimethyl Phthalate	1.396	1.425	-2.1	111	0.00
44	TM	Acenaphthylene	1.860	1.872	-0.6	107	0.00
45	TM	3-Nitroaniline	0.397	0.412	-3.8	109	0.00
46	TMC	Acenaphthene	1.057	1.069	-1.1	111	0.00
47	TPM	2,4-dinitrophenol	0.167	0.198	-18.6	115	0.00
48	TM	Dibenzofuran	1.661	1.682	-1.3	111	0.00
49	TMP	4-Nitrophenol	0.165	0.172	-4.2	121	0.00
50	TM	2,4-Dinitrotoluene	0.450	0.466	-3.6	108	0.00
51	T	2,3,4,6-Tetrachlorophenol	0.250	0.272	-8.8	115	0.00
52	TM	Fluorene	1.296	1.283	1.0	108	0.00
	TM	Diethyl Phthalate	1.380	1.387	-0.5	107	0.00
54	TM	4-Chlorophenyl phenyl ethe	0.610	0.622	-2.0	111	0.00
55	TM	4-Nitroaniline	0.403	0.422	-4.7	107	0.00
56	TM	4,6-Dinitro-2-methylphenol	0.266	0.300	-12.8	111	0.00
57	TCM	n-nitrosodiphenylamine	0.876	0.870	0.7	108	0.00
58	TM	Azobenzene	1.897	1.828	3.6	103	0.00
59	TM	Biphenyl	1.513	1.504	0.6	111	0.00
60	S	2,4,6-Tribromophenol	0.214	0.224	-4.7	114	0.00
61	I	d10-Phenanthrene	1.000	1.000	0.0	108	0.00
62	TM	4-bromophenyl phenyl ether	0.216	0.220	-1.9	110	0.00
63	TM	Hexachlorobenzene	0.251	0.254	-1.2	105	0.00
64	TMC	Pentachlorophenol	0.117	0.136	-16.2	131	0.00
65	TM	Phenanthrene	1.009	1.013	-0.4	109	0.00
66	TM	Anthracene	1.018	1.016	0.2	104	0.00
67	TM	Carbazole	0.984	1.015	-3.2	110	0.00
68	TM	Di-n-butyl phthalate	1.367	1.366	0.1	102	0.00
69	TCM	Fluoranthene	1.039	1.134	-9.1	115	0.00
70	TM	Benzidine	0.507	0.124	75.5#	25#	0.00
71	I	d12-Chrysene	1.000	1.000	0.0	104	0.00
72	TM	Pyrene	1.204	1.372	-14.0	118	0.00
73	S	d14-p-terphenyl	0.855	0.903	-5.6	111	0.00
74	TM	Butyl benzyl phthalate	0.717	0.742	-3.5	106	0.00
75	TM	Benzo[a]anthracene	1.071	1.088	-1.6	105	0.00
76	TM	Chrysene	1.011	1.034	-2.3	105	0.00
	TM	3,3'-Dichlorobenzidine	0.455	0.450	1.1	103	0.00
78	TM	Bis (2-ethylhexyl) phthalat	1.000	0.984	1.6	103	0.00

(#) = Out of Range

Evaluate Continuing Calibration Report

Data File : D:\HPCHEM\1\DATA\053106-D\D47795SC.D Vial: 2
Acq On : 31 May 2006 11:03 am Operator:
Sample : S6849 PT3 20/40 Inst : D Instrum
Misc : Multiplr: 1.00
MS Integration Params: RTEINT.P

Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Wed May 31 12:09:53 2006
Response via : Multiple Level Calibration

Min. RRF : 0.050 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
Max. RRF Dev : 20% Max. Rel. Area : 200%

		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
79	I	d12-Perylene	1.000	1.000	0.0	95	0.00
80	TCM	di-n-octyl-phthalate	1.764	1.860	-5.4	98	0.00
81	TM	Benzo[b] fluoranthene	1.103	1.100	0.3	94	0.00
82	TM	Benzo[k] fluoranthene	1.063	1.109	-4.3	97	0.00
83	TCM	Benzo[a] pyrene	1.042	1.011	3.0	91	0.00
84	TM	Indeno [1,2,3-cd] pyrene	1.023	1.026	-0.3	96	0.00
85	TM	Dibenz [a,h] anthracene	0.998	0.971	2.7	92	0.00
86	TM	Benzo(g,h,i) perylene	1.002	0.979	2.3	92	0.00

(#) = Out of Range SPCC's out = 0 CCC's out = 0

Quantitation Report (Not Reviewed)

Data File : D:\HPCHEM\1\DATA\053106-D\D47796SC.D Vial: 3
Acq On : 31 May 2006 11:40 am Operator:
Sample : S6740 B+A 20 Inst : D Instrum
Misc : Multiplr: 1.00
IS Integration Params: RTEINT.P
Quant Time: May 31 12:10 2006 Quant Results File: S05246.RES

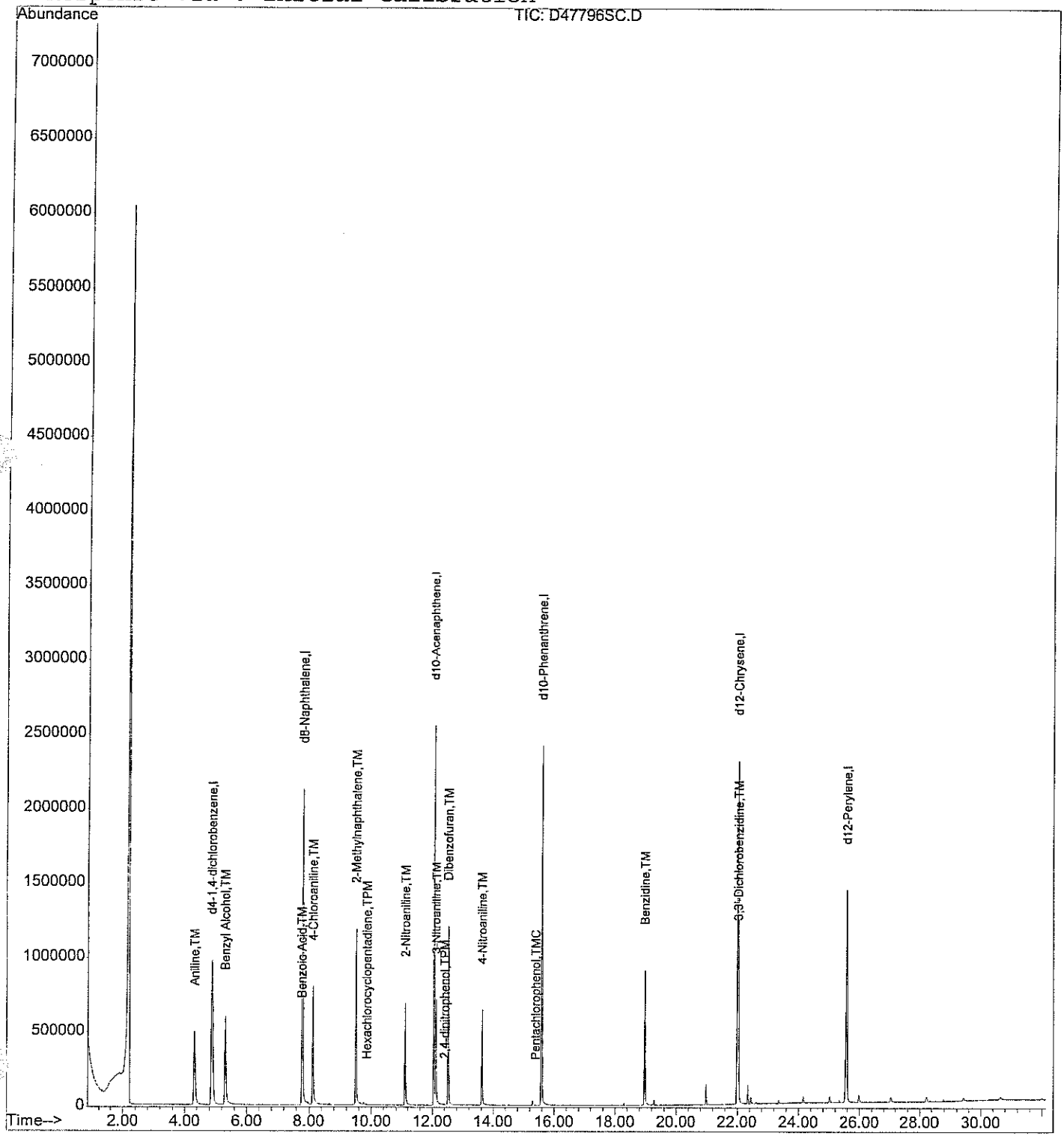
Quant Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Wed May 31 12:09:53 2006
Response via : Initial Calibration
DataAcq Meth : ABN

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) d4-1,4-dichlorobenzene	4.88	152	563613	40.00	ug/ml	0.00
19) d8-Naphthalene	7.79	136	2115621	40.00	ug/ml	0.00
35) d10-Acenaphthene	12.05	164	1041300	40.00	ug/ml	0.00
61) d10-Phenanthrene	15.59	188	1804982	40.00	ug/ml	0.00
71) d12-Chrysene	22.01	240	1533577	40.00	ug/ml	0.00
79) d12-Perylene	25.58	264	1372239	40.00	ug/ml	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	2.57	112	108	0.00	ug/ml	0.00
Spiked Amount 75.000	Range 28 - 110		Recovery =	0.00%#		
5) d5-Phenol	4.42	99	177	0.01	ug/ml	-0.01
Spiked Amount 75.000	Range 22 - 110		Recovery =	0.01%#		
20) d5-nitrobenzene	6.24	82	50	0.00	ug/ml	0.02
Spiked Amount 50.000	Range 45 - 107		Recovery =	0.00%#		
39) 2-Fluorobiphenyl	10.50	172	46	0.00	ug/ml	0.00
Spiked Amount 50.000	Range 44 - 102		Recovery =	0.00%#		
60) 2,4,6-Tribromophenol	13.89	330	43	0.01	ug/ml	-0.13
Spiked Amount 75.000	Range 63 - 127		Recovery =	0.01%#		
73) d14-p-terphenyl	19.56	244	62	0.00	ug/ml	0.00
Spiked Amount 50.000	Range 33 - 132		Recovery =	0.00%#		
Target Compounds						Qvalue
6) Aniline	4.32	93	674693	19.64	ug/ml	98
13) Benzyl Alcohol	5.31	108	287037	19.29	ug/ml	97
27) Benzoic Acid	7.76	105	154	5.18	ug/ml#	1
31) 4-Chloroaniline	8.12	127	491322	19.02	ug/ml	95
34) 2-Methylnaphthalene	9.50	141	576700	17.22	ug/ml#	58
36) Hexachlorocyclopentadiene	9.85	237	85	5.49	ug/ml#	7
41) 2-Nitroaniline	11.11	138	264683	20.24	ug/ml	98
45) 3-Nitroaniline	12.13	138	215817	20.86	ug/ml	94
47) 2,4-dinitrophenol	12.41	184	67	5.39	ug/ml#	39
48) Dibenzofuran	12.52	168	841653	19.47	ug/ml	100
55) 4-Nitroaniline	13.63	138	216153	20.61	ug/ml	98
64) Pentachlorophenol	15.39	266	45	6.50	ug/ml#	17
70) Benzidine	18.97	184	641059	28.01	ug/ml	99
77) 3,3'-Dichlorobenzidine	22.03	252	309137	17.74	ug/ml	98

(#) = qualifier out of range (m) = manual integration

Data File : D:\HPCHEM\1\DATA\053106-D\D47796SC.D Vial: 3
Acq On : 31 May 2006 11:40 am Operator:
Sample : S6740 B+A 20 Inst : D Instrum
Misc : Multiplr: 1.00
MS Integration Params: RTEINT.P
Quant Time: May 31 12:10 2006 Quant Results File: S05246.RES

Method : D:\HPCHEM\1\METHODS\S05246.M (RTE Integrator)
Title : ABN
Last Update : Wed May 31 12:09:53 2006
Response via : Initial Calibration



SEMI-VOLATILE ORGANIC
INTERNAL STANDARD AREA AND RT SUMMARY

Project #: 218736.002 | Quote #: NA | SDG#: 56637

DFTPP File ID: D47722T.D | Site ID: GP Old Town Mill | Location: ME

Instrument ID: D | DFTPP Injection Date: 5/25/06

GC Column: RTX-5MS | ID: 0.25(mm) | DFTPP Injection Time: 0305

	IS1 AREA #	RT #	IS2 AREA #	RT #	IS3 AREA #	RT #	IS4 AREA #	RT #	IS5 AREA #	RT #	IS6 AREA #	RT #
12 HOUR STD	487489	5.04	1805490	7.95	924277	12.21	1566340	15.75	1389790	22.17	1317030	25.81
UPPER LIMIT	974978	5.34	3610980	8.43	1848554	12.95	3132680	16.69	2779580	23.50	2634060	27.36
LOWER LIMIT	243745	4.74	902745	7.48	462139	11.48	783170	14.80	694895	20.84	658515	24.26
SAMPLE NO.												
01 S6852 PT1 2/4	499606	5.04	1884970	7.95	940184	12.21	1589690	15.75	1403450	22.17	1349550	25.80
02 S6851 PT2A 4/8	495838	5.03	1830300	7.95	941311	12.21	1609030	15.75	1433200	22.16	1362420	25.80
03 S6850 PT2 10/20	492927	5.04	1801960	7.95	930235	12.21	1577640	15.75	1404560	22.17	1346260	25.80
04 S6849 PT3 20/40	487489	5.04	1805490	7.95	924277	12.21	1566340	15.75	1389790	22.17	1317030	25.81
05 S6848 PT4A 40/80	494381	5.04	1825430	7.96	917039	12.22	1549650	15.76	1392900	22.18	1308520	25.81
06 S6847 PT4 50/100	486916	5.04	1808660	7.96	934498	12.22	1556870	15.76	1376000	22.17	1311670	25.82
07 S6846 PT5 100/200	502892	5.05	1847600	7.98	942480	12.23	1557200	15.77	1196580	22.19	1227070	25.83
08 S6857 ABN CC 20/40	491009	5.04	1822320	7.95	917674	12.21	1589010	15.75	1367670	22.17	1331050	25.80
09 S6858 BENZIDINES + ANILI	481737	5.03	1820100	7.95	905390	12.21	1525850	15.74	1385820	22.16	1359280	25.80
10												
11												
12												
13												
14												
15												
16												
17												
18												
19												
20												
21												
22												

IS1 = d4-1,4 -Dichlorobenzene
IS2 =d8- Naphthalene
IS3 = d10-Acenaphthalene
IS4 = d10- Phenanthrene
IS5 = d12-Chrysene
IS6 = d10-Perylene

AREA UPPER LIMIT = +100 of internal standard area
AREA LOWER LIMIT = -50 of internal standard area
RT UPPER LIMIT = +0.06 minutes of internal standard RT
RT LOWER LIMIT = -0.06 minutes of internal standard RT

Column used to flag values outside QC limits with an asterisk
* Values outside of QC limits.

368/486

SEMI-VOLATILE ORGANIC
INTERNAL STANDARD AREA AND RT SUMMARY

Project #: 218736.002 | Quote #: NA | SDG#: 56637

DFTPP File ID: D47755T.D | Site ID: GP Old Town Mill | Location: ME

Instrument ID: D | DFTPP Injection Date: 5/30/06

GC Column: RTX-5MS | ID: 0.25(mm) | DFTPP Injection Time: 0639

	IS1 AREA #	RT #	IS2 AREA #	RT #	IS3 AREA #	RT #	IS4 AREA #	RT #	IS5 AREA #	RT #	IS6 AREA #	RT #
12 HOUR STD	582194	4.98	2153580	7.90	1146580	12.16	2037120	15.70	1830140	22.12	1910720	25.75
UPPER LIMIT	1164388	5.28	4307160	8.37	2293160	12.89	4074240	16.64	3660280	23.45	3821440	27.29
LOWER LIMIT	291097	4.69	1076790	7.42	573290	11.43	1018560	14.76	915070	20.80	955360	24.20
SAMPLE NO.												
01 S6849 PT3 20/40	582194	4.98	2153580	7.90	1146580	12.16	2037120	15.70	1830140	22.12	1910720	25.75
02 S6760 B+A 20	620189	4.98	2352120	7.89	1202450	12.15	2170250	15.69	1951560	22.11	2018810	25.73
03 B05306AW	607093	4.98	2235920	7.89	1145160	12.15	2031610	15.70	1820290	22.11	1823140	25.73
04 L05306AW	582705	4.99	2165830	7.90	1136000	12.16	1972550	15.69	1762420	22.12	1815750	25.74
05 LD05306AW	623379	4.98	2284570	7.89	1192280	12.16	2069180	15.70	1813950	22.12	1801020	25.74
06 56637-1	599539	4.98	2217420	7.89	1144900	12.15	1993730	15.69	1634620	22.12	1825950	25.75
07 56637-2	610509	4.98	2327820	7.89	1178580	12.15	2070450	15.69	1716850	22.11	1740760	25.73
08 56637-3	573225	4.98	2161080	7.89	1086340	12.15	1930690	15.69	1646830	22.11	1656180	25.73
09 56637-20	560338	4.98	2134060	7.89	1036140	12.15	1812030	15.70	1568660	22.11	1693300	25.74
10 56637-21	566110	4.98	2159890	7.89	1091400	12.15	1962840	15.69	1664050	22.11	1659260	25.73
11 56637-1, RE-INJECT	597641	4.98	2188400	7.89	1134090	12.15	1976180	15.69	1657130	22.11	1849000	25.74
12 56637-2, RE-INJECT	605541	4.98	2266930	7.89	1145300	12.15	2033040	15.69	1702580	22.11	1710780	25.73
13 56637-20, RE-INJE	557049	4.98	2100140	7.89	1045230	12.15	1836110	15.69	1544570	22.11	1674660	25.74
14 56637-21, RE-INJE	570466	4.98	2150250	7.89	1087150	12.15	1951500	15.69	1606080	22.11	1632160	25.73
15 LD05306AW, RE-INJECT	602531	4.99	2223300	7.90	1152220	12.16	1999300	15.70	1714120	22.12	1674840	25.74
16												
17												
18												
19												
20												
21												
22												

IS1 = d4-1,4 -Dichlorobenzene
IS2 =d8- Naphthalene
IS3 = d10-Acenaphthalene
IS4 = d10- Phenanthrene
IS5 = d12-Chrysene
IS6 = d10-Perylene

AREA UPPER LIMIT = +100 of internal standard area
AREA LOWER LIMIT = -50 of internal standard area
RT UPPER LIMIT = +0.06 minutes of internal standard RT
RT LOWER LIMIT = -0.06 minutes of internal standard RT

Column used to flag values outside QC limits with an asterisk
* Values outside of QC limits.

SEMI-VOLATILE ORGANIC
INTERNAL STANDARD AREA AND RT SUMMARY

Project #: 218736.002 | Quote #: NA | SDG#: 56637

DFTPP File ID: D47771T.D | Site ID: GP Old Town Mill | Location: ME

Instrument ID: D | DFTPP Injection Date: 5/30/06

GC Column: RTX-5MS | ID: 0.25(mm) | DFTPP Injection Time: 2043

	IS1 AREA	#	RT	#	IS2 AREA	#	RT	#	IS3 AREA	#	RT	#	IS4 AREA	#	RT	#	IS5 AREA	#	RT	#	IS6 AREA	#	RT	#
12 HOUR STD	595490		4.97		2193200		7.89		1141720		12.15		1972470		15.70		1676490		22.11		1552180		25.72	
UPPER LIMIT	1190980		5.27		4386400		8.37		2283440		12.88		3944940		16.64		3352980		23.44		3104360		27.26	
LOWER LIMIT	297745		4.67		1096600		7.42		570860		11.42		986235		14.75		838245		20.78		776090		24.18	
SAMPLE NO.																								
01 S6849 PT3 20/40	595490		4.97		2193200		7.89		1141720		12.15		1972470		15.70		1676490		22.11		1552180		25.72	
02 S6760 B+A 20	631919		4.97		2372910		7.88		1211650		12.14		2126590		15.69		1795090		22.10		1664860		25.71	
03 B05306AS	613418		4.98		2325330		7.88		1171480		12.14		2023830		15.69		1688560		22.10		1551940		25.71	
04 L05306AS	635428		4.98		2359190		7.89		1203980		12.15		2098360		15.69		1618440		22.11		1394920		25.72	
05 LD05306AS	624028		4.98		2293180		7.89		1171430		12.15		2021440		15.69		1695790		22.11		1531200		25.72	
06 ZZZZZ	629067		4.97		2326210		7.88		1180140		12.14		2011360		15.68		1552270		22.10		1518320		25.72	
07 ZZZZZ	643396		4.97		2422410		7.88		1211070		12.14		2124510		15.68		1618230		22.11		1652540		25.72	
08 56637-4	634562		4.97		2389050		7.88		1223980		12.14		2093370		15.69		1624330		22.10		1536550		25.72	
09 56637-9	625739		4.97		2330250		7.88		1160800		12.15		2061990		15.68		1682190		22.10		1556360		25.71	
10 56637-10	630508		4.97		2395770		7.88		1194410		12.14		2067650		15.69		1672220		22.10		1535800		25.72	
11 56637-12	613629		4.97		2306200		7.89		1162130		12.15		2017900		15.68		1653670		22.10		1490750		25.71	
12 56637-18	626664		4.97		2261420		7.89		1233520		12.16		2516400		15.72		1501810		22.11		1521010		25.72	
13 56637-11	608781		4.97		2254830		7.89		1144590		12.15		2026440		15.68		1629500		22.11		1555800		25.72	
14 56637-14	587998		4.97		2191320		7.88		1094820		12.14		1828590		15.69		2004580		22.18		318971	*	25.81	
15 56637-19	559898		4.98		2100210		7.89		1071140		12.15		1848020		15.69		1890440		22.11		1227280		25.72	
16 56637-22	580618		4.98		2208030		7.88		1120270		12.15		1881590		15.70		1572460		22.12		421984	*	25.73	
17 56637-23	584683		4.97		2209060		7.89		1100840		12.1476		1857640		15.69		1502400		22.12		415856	*	25.72	
18 56637-15,,1:2	550858		4.98		2136070		7.89		1097540		12.1767		2569930		15.82		984808		22.26		146891	*	25.78	
19 56637-16,,1:5	558887		4.98		2066570		7.90		1300600		12.1763		2509880		15.78		1183650		22.19		200895	*	25.76	
20 56637-17,,1:5	560323		4.98		2032870		7.89		1036530		12.1519		1893820		15.71		818343	*	22.16		176058	*	25.75	
21																								
22																								

IS1 = d4-1,4 -Dichlorobenzene
IS2 =d8- Naphthalene
IS3 = d10-Acenaphthalene
IS4 = d10- Phenanthrene
IS5 = d12-Chrysene
IS6 = d10-Perylene

AREA UPPER LIMIT = +100 of internal standard area
AREA LOWER LIMIT = -50 of internal standard area
RT UPPER LIMIT = +0.06 minutes of internal standard RT
RT LOWER LIMIT = -0.06 minutes of internal standard RT

Column used to flag values outside QC limits with an asterisk
* Values outside of QC limits.

370/486

SEMI-VOLATILE ORGANIC
INTERNAL STANDARD AREA AND RT SUMMARY

Project #: 218736.002 | Quote #: NA | SDG#: 56637

DFTPP File ID: D47794T.D | Site ID: GP Old Town Mill | Location: ME

Instrument ID: D | DFTPP Injection Date: 5/31/06

GC Column: RTX-5MS | ID: 0.25(mm) | DFTPP Injection Time: 1018

	IS1 AREA #	RT #	IS2 AREA #	RT #	IS3 AREA #	RT #	IS4 AREA #	RT #	IS5 AREA #	RT #	IS6 AREA #	RT #
12 HOUR STD	546320	4.88	1990730	7.80	1010440	12.06	1685500	15.60	1450960	22.01	1254130	25.58
UPPER LIMIT	1092640	5.17	3981460	8.27	2020880	12.79	3371000	16.53	2901920	23.33	2508260	27.12
LOWER LIMIT	273160	4.59	995365	7.33	505220	11.34	842750	14.66	725480	20.69	627065	24.05
SAMPLE NO.												
01 S6849 PT3 20/40	546320	4.88	1990730	7.80	1010440	12.06	1685500	15.60	1450960	22.01	1254130	25.58
02 S6740 B+A 20	563613	4.88	2115620	7.79	1041300	12.05	1804980	15.59	1533580	22.01	1372240	25.58
03 B05306AS	524660	4.88	1978960	7.79	984615	12.05	1701440	15.59	1406590	22.00	1235340	25.57
04 56637-15,,1:100	481269	4.89	1818230	7.79	904598	12.05	1512860	15.59	1260660	22.01	1113830	25.59
05 56637-16,,1:50	496119	4.89	1825190	7.80	881206	12.06	1465750	15.60	1306470	22.02	1285790	25.60
06 56637-17,,1:100	491814	4.88	1827130	7.79	897434	12.05	1497930	15.59	1282340	22.01	1212020	25.59
07 56637-14,DL,1:2	512188	4.88	1906280	7.79	917852	12.05	1678650	15.60	1737610	22.06	1535630	25.72
08 56637-22,DL,1:2	489601	4.89	1823890	7.80	905103	12.06	1530530	15.60	1532790	22.03	1164200	25.62
09 56637-23,DL,1:2	501481	4.89	1861660	7.80	931608	12.06	1667160	15.60	1442120	22.03	1252150	25.61
10 56637-17,,1:20	455954	4.89	1726010	7.80	859385	12.06	1508160	15.61	1427740	22.04	861526	25.63
11 56637-23,MS,,1:2	477888	4.89	1777510	7.80	898618	12.06	1599840	15.61	1435210	22.03	795076	25.61
12 56637-23,MSD,,1:2	496469	4.90	1833110	7.81	923808	12.07	1649480	15.61	1486540	22.03	806253	25.62
13 ZZZZZ	506459	4.90	1901350	7.81	921329	12.06	1708550	15.61	1515860	22.03	778805	25.60
14 ZZZZZ	489542	4.90	1825710	7.80	899472	12.07	1620350	15.61	1493010	22.03	844719	25.61
15 ZZZZZ	506485	4.90	1884060	7.82	1112890	12.10	3016290	15.67	1667620	22.15	384752*	25.73
16 ZZZZZ	500862	4.90	1871180	7.81	973611	12.07	1915270	15.62	1793740	22.05	483240*	25.62
17												
18												
19												
20												
21												
22												

IS1 = d4-1,4 -Dichlorobenzene
IS2 =d8- Naphthalene
IS3 = d10-Acenaphthalene
IS4 = d10- Phenanthrene
IS5 = d12-Chrysene
IS6 = d10-Perylene

AREA UPPER LIMIT = +100 of internal standard area
AREA LOWER LIMIT = -50 of internal standard area
RT UPPER LIMIT = +0.06 minutes of internal standard RT
RT LOWER LIMIT = -0.06 minutes of internal standard RT

Column used to flag values outside QC limits with an asterisk
* Values outside of QC limits.

371486

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EXTRACTION
LOGBOOK PAGES

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PCB
DATA SUMMARIES

analytics environmental laboratory LLC

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800-929-9906

Mr. Karl Kasper
Woodard & Curran
41 Hutchins Drive
Portland ME 04102

June 6, 2006

CLIENT SAMPLE ID

Project Name: GP Old Town Mill

Project Number: 218736.002

Field Sample ID: LABQC

SAMPLE DATA

Lab Sample ID: B05306PAS
Matrix: Solid
Percent Solid: 100
Dilution Factor: 1.0
Collection Date:
Lab Receipt Date:
Extraction Date: 05/30/06
Analysis Date: 05/30/06

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit µg/kg	Results µg/kg
PCB-1016	15	U
PCB-1221	15	U
PCB-1232	15	U
PCB-1242	15	U
PCB-1248	15	U
PCB-1254	15	U
PCB-1260	15	U

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	64	%
Decachlorobiphenyl	44	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082.

Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3545.

COMMENTS: Results are expressed on a dry weight basis.

PCB Report

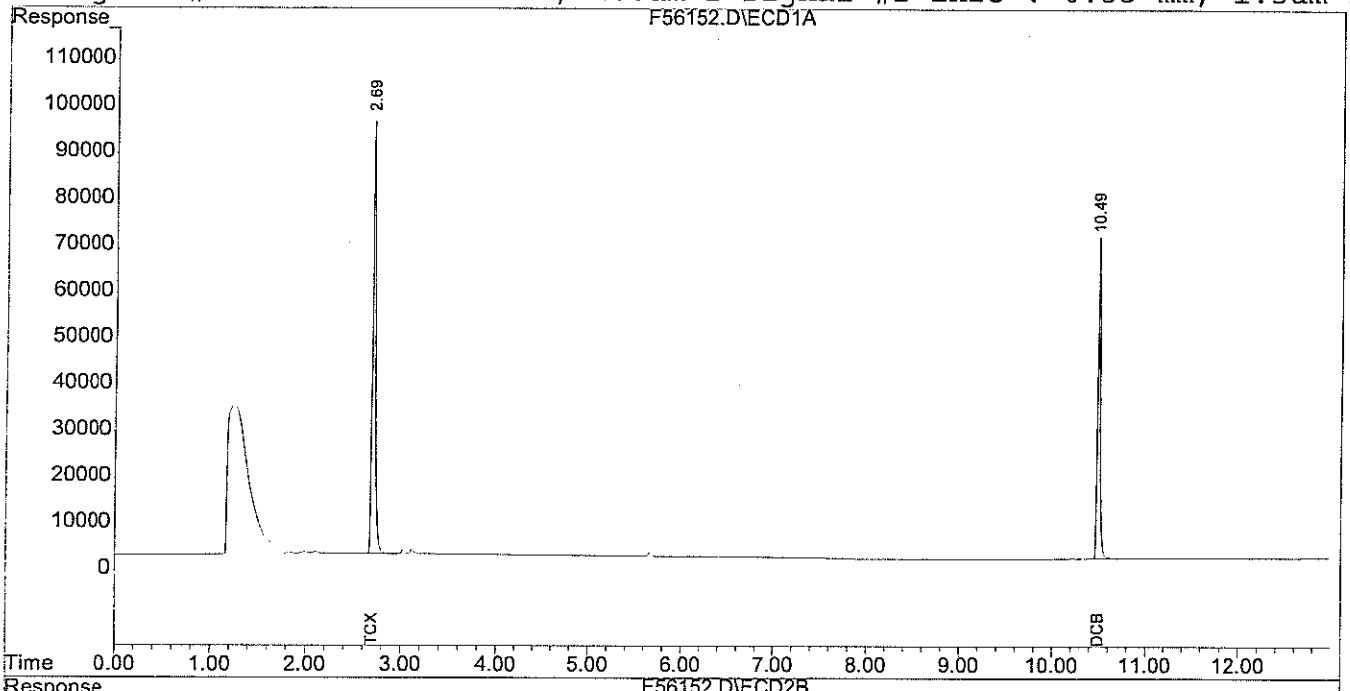
Authorized signature

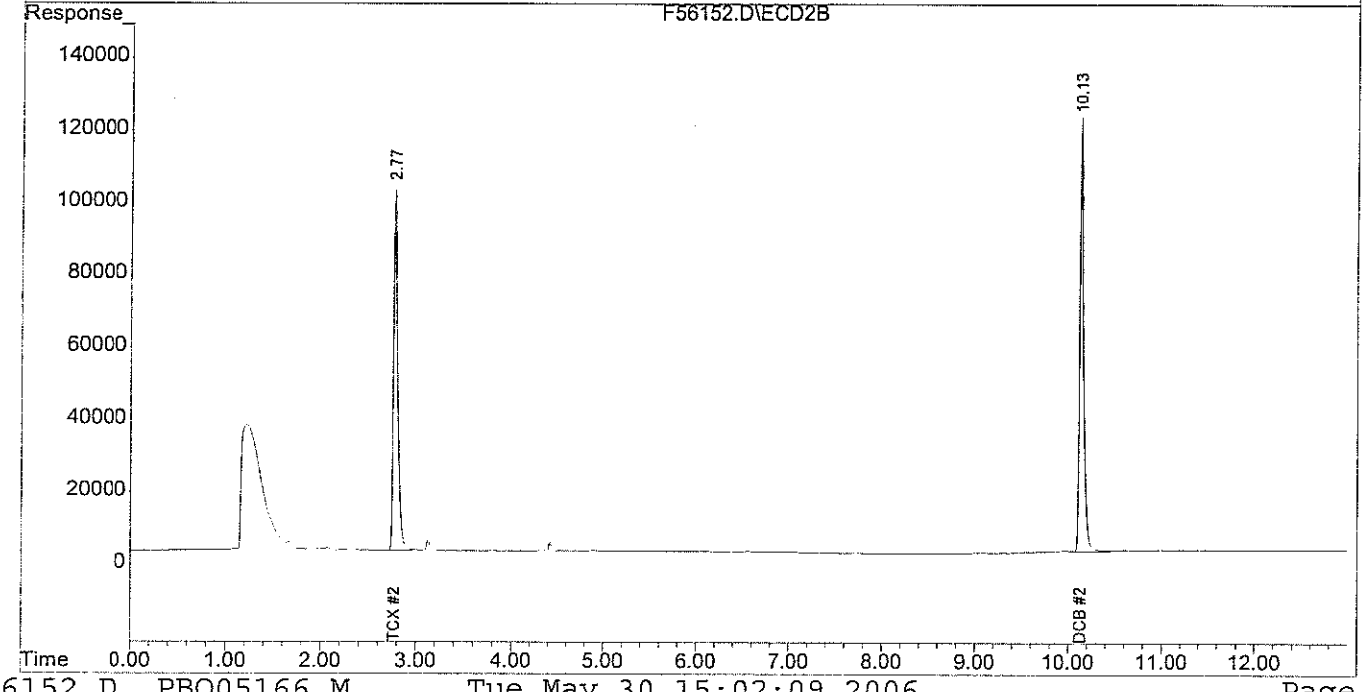
Quantitation Report (QT Reviewed)

Signal #1 : C:\HPCHEM\2\DATA\053006-F\F56152.D\ECD1A.CH Vial: 5
Signal #2 : C:\HPCHEM\2\DATA\053006-F\F56152.D\ECD2B.CH
Acq On : 30 May 2006 2:27 pm Operator:
Sample : B05306PAS, A/C Inst : Instrumen
Misc : SOIL Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 30 14:58 2006 Quant Results File: PBQ05166.RES

Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 10:06:32 2006
Response via : Multiple Level Calibration
DataAcq Meth : PSTQUICK.M

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film





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Mr. Karl Kasper
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41 Hutchins Drive
Portland ME 04102

June 6, 2006

CLIENT SAMPLE ID

Project Name: GP Old Town Mill

Project Number: 218736.002

Field Sample ID: LABQC

SAMPLE DATA

Lab Sample ID: B05306PAS RR
Matrix: Solid
Percent Solid: 100
Dilution Factor: 1.0
Collection Date:
Lab Receipt Date:
Extraction Date: 05/30/06
Analysis Date: 05/31/06

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit µg/kg	Results µg/kg
PCB-1016	15	U
PCB-1221	15	U
PCB-1232	15	U
PCB-1242	15	U
PCB-1248	15	U
PCB-1254	15	U
PCB-1260	15	U

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	88	%
Decachlorobiphenyl	92	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082.

Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3545.

COMMENTS: Results are expressed on a dry weight basis.

PCB Report

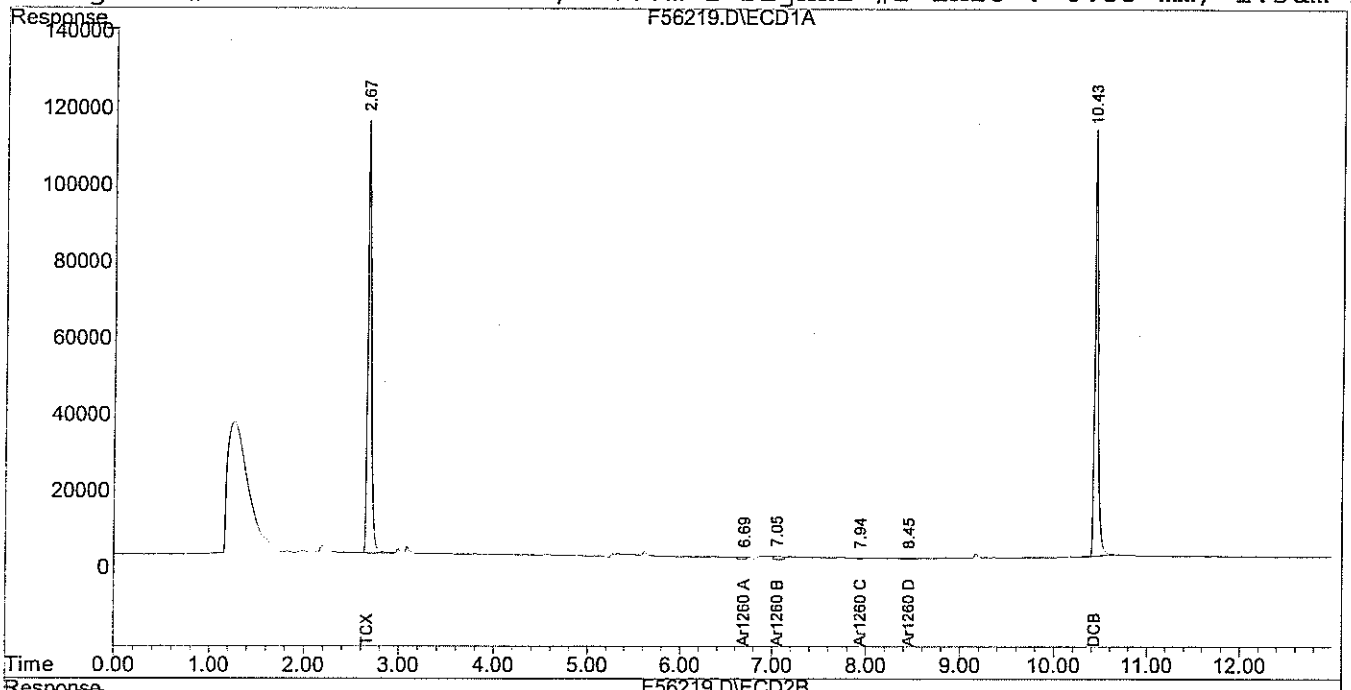
Authorized signature

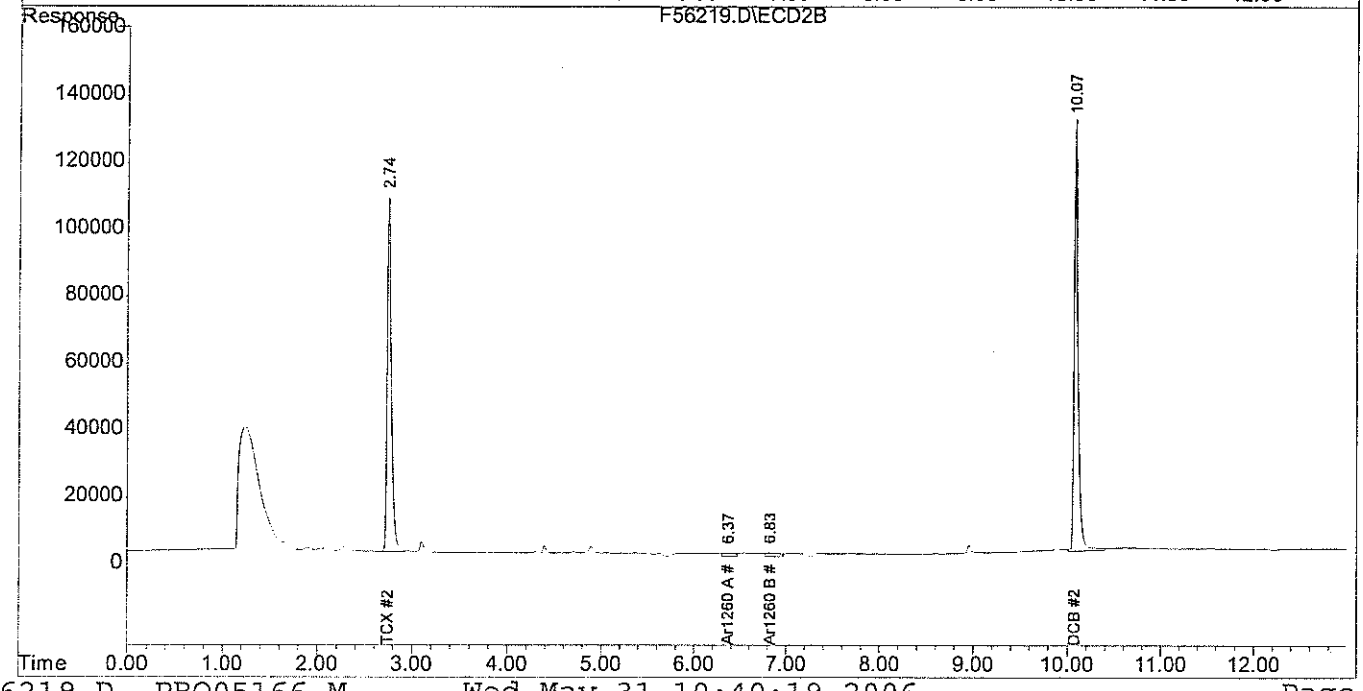
Quantitation Report (Not Reviewed)

Signal #1 : C:\HPCHEM\2\DATA\053106-F\F56219.D\ECD1A.CH Vial: 2
Signal #2 : C:\HPCHEM\2\DATA\053106-F\F56219.D\ECD2B.CH
Acq On : 31 May 2006 9:58 am Operator:
Sample : B05306PAS, A/C, RR Inst : Instrumen
Misc : SOIL Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 31 10:40 2006 Quant Results File: PBQ05166.RES

Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 10:06:32 2006
Response via : Multiple Level Calibration
DataAcq Meth : PSTQUICK.M

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film





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Portland ME 04102

May 31, 2006

CLIENT SAMPLE ID

Project Name: GP Old Town Mill

Project Number: 218736.002

Field Sample ID: SB-22 (0-4)

SAMPLE DATA

Lab Sample ID: 56637-4
Matrix: Solid
Percent Solid: 85
Dilution Factor: 1.2
Collection Date: 05/25/06
Lab Receipt Date: 05/27/06
Extraction Date: 05/30/06
Analysis Date: 05/30/06

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit µg/kg	Results µg/kg
PCB-1016	18	U
PCB-1221	18	U
PCB-1232	18	U
PCB-1242	18	U
PCB-1248	18	U
PCB-1254	18	U
PCB-1260	18	U

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	92	%
Decachlorobiphenyl	81	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082.

Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3545.

COMMENTS: Results are expressed on a dry weight basis.

PCB Report

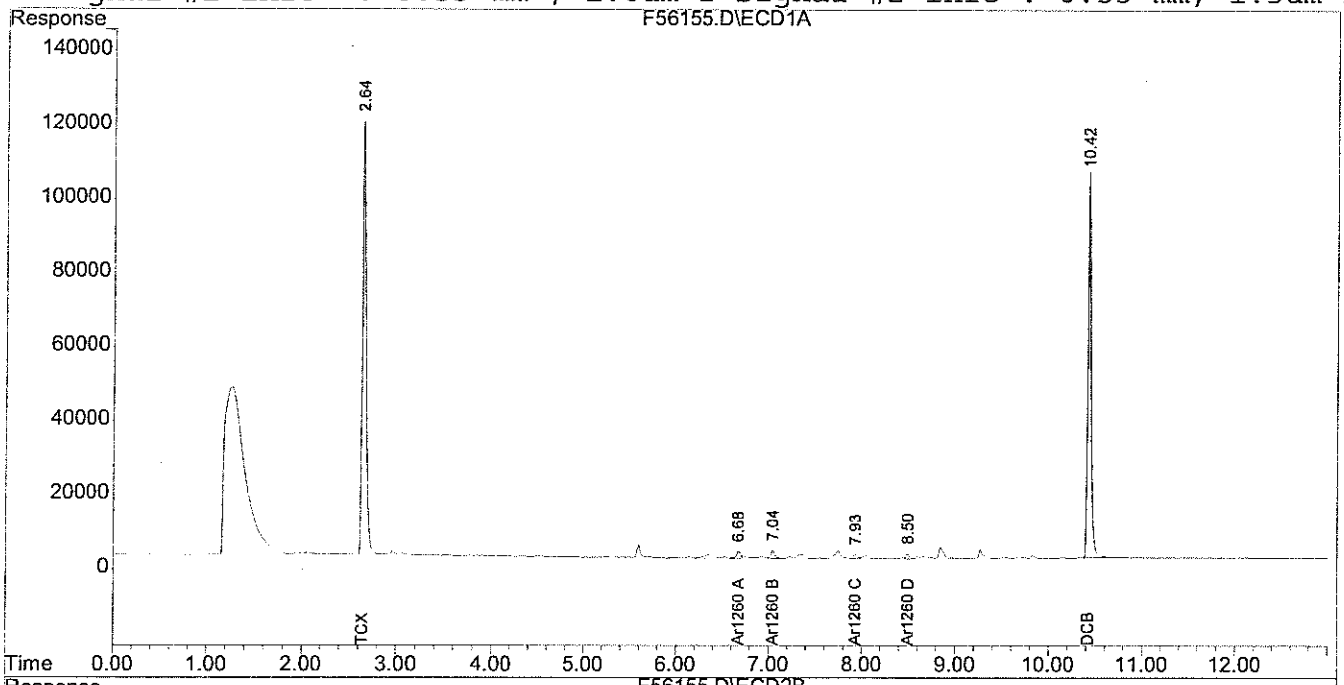
Authorized signature Melissa Tall

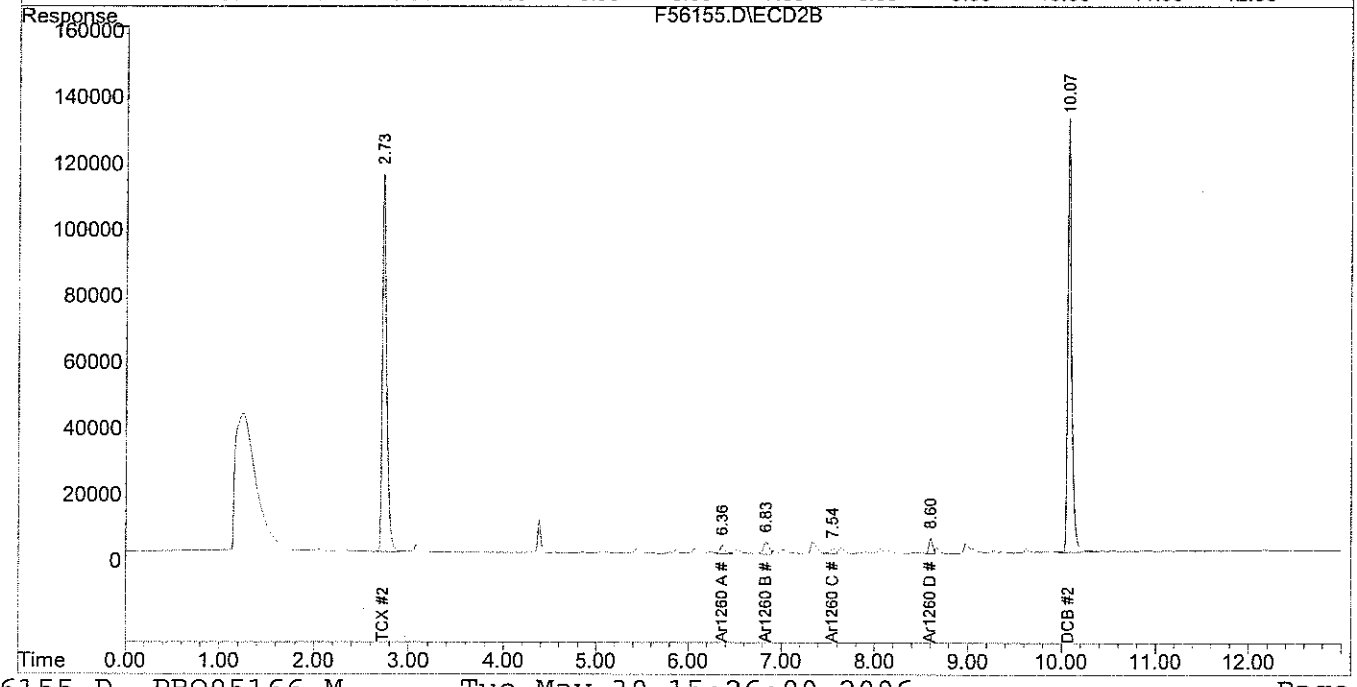
Quantitation Report (QT Reviewed)

Signal #1 : C:\HPCHEM\2\DATA\053006-F\F56155.D\ECD1A.CH Vial: 8
Signal #2 : C:\HPCHEM\2\DATA\053006-F\F56155.D\ECD2B.CH
Acq On : 30 May 2006 3:11 pm Operator:
Sample : 56637-4, A/C Inst : Instrumen
Misc : SOIL Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 30 15:25 2006 Quant Results File: PBQ05166.RES

Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 10:06:32 2006
Response via : Multiple Level Calibration
DataAcq Meth : PSTQUICK.M

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film





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Mr. Karl Kasper
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41 Hutchins Drive
Portland ME 04102

May 31, 2006

CLIENT SAMPLE ID

Project Name: GP Old Town Mill

Project Number: 218736.002

Field Sample ID: SB-21 (4-8')

SAMPLE DATA

Lab Sample ID: 56637-6
Matrix: Solid
Percent Solid: 95
Dilution Factor: 1.0
Collection Date: 05/25/06
Lab Receipt Date: 05/27/06
Extraction Date: 05/30/06
Analysis Date: 05/30/06

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit µg/kg	Results µg/kg
PCB-1016	15	U
PCB-1221	15	U
PCB-1232	15	U
PCB-1242	15	U
PCB-1248	15	U
PCB-1254	15	**19 D**
PCB-1260	15	U

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	98	%
Decachlorobiphenyl	94	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082.

Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3545.

COMMENTS: Results are expressed on a dry weight basis. D=Sample did not meet confirmation acceptance criteria for percent difference.

PCB Report

Authorized signature

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PCB CONFIRMATION RESULTS

SDG: 56637

Project No: 218736.002 Site ID: Location: ME

Instrument ID: F GC Columns: RTX-CLPesticidesI ID: 0.32(mm)

RTX-CLPesticidesII

Sample ID: 56637-6

COMPOUND	RESPONSE #1 ug/kg	RESPONSE #2 ug/kg	DIFFERENCE %	#	% DIFFERENCE Acceptance Limit
PCB 1254	12 J	19	47	*	40

Column to be used to flag % Difference with an asterisk

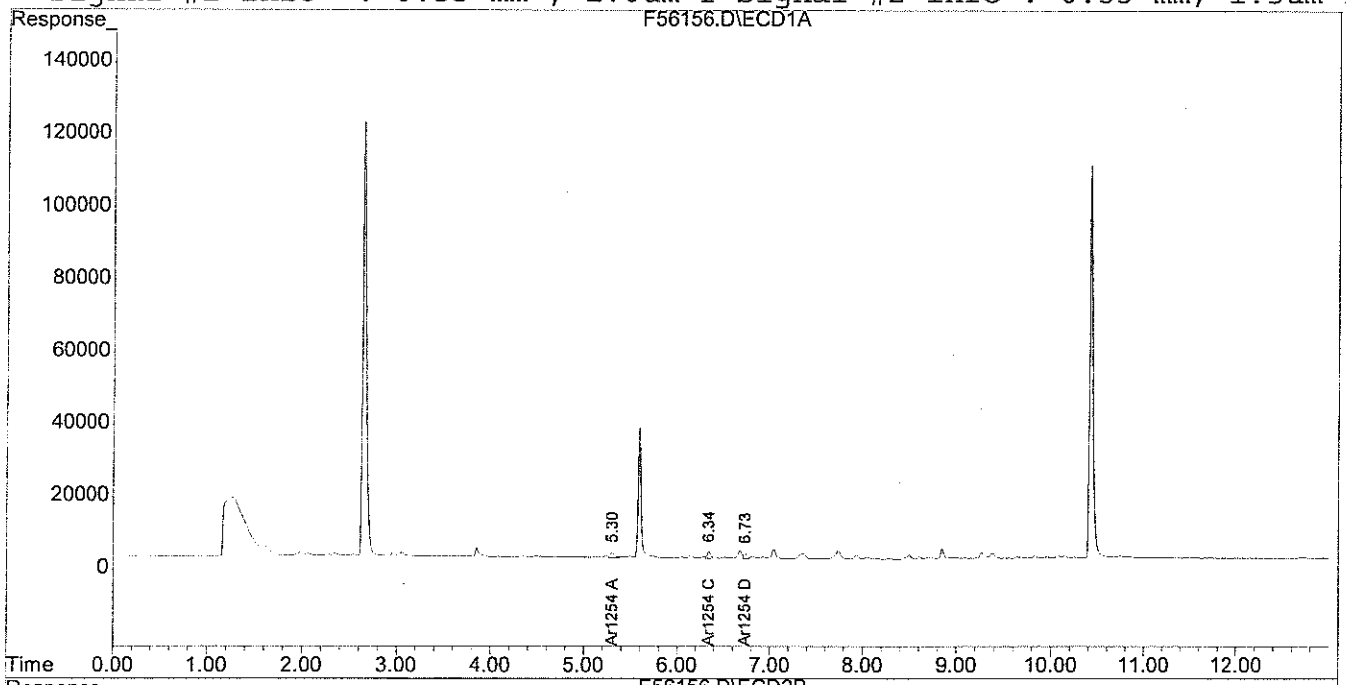
* Values outside of QC limits

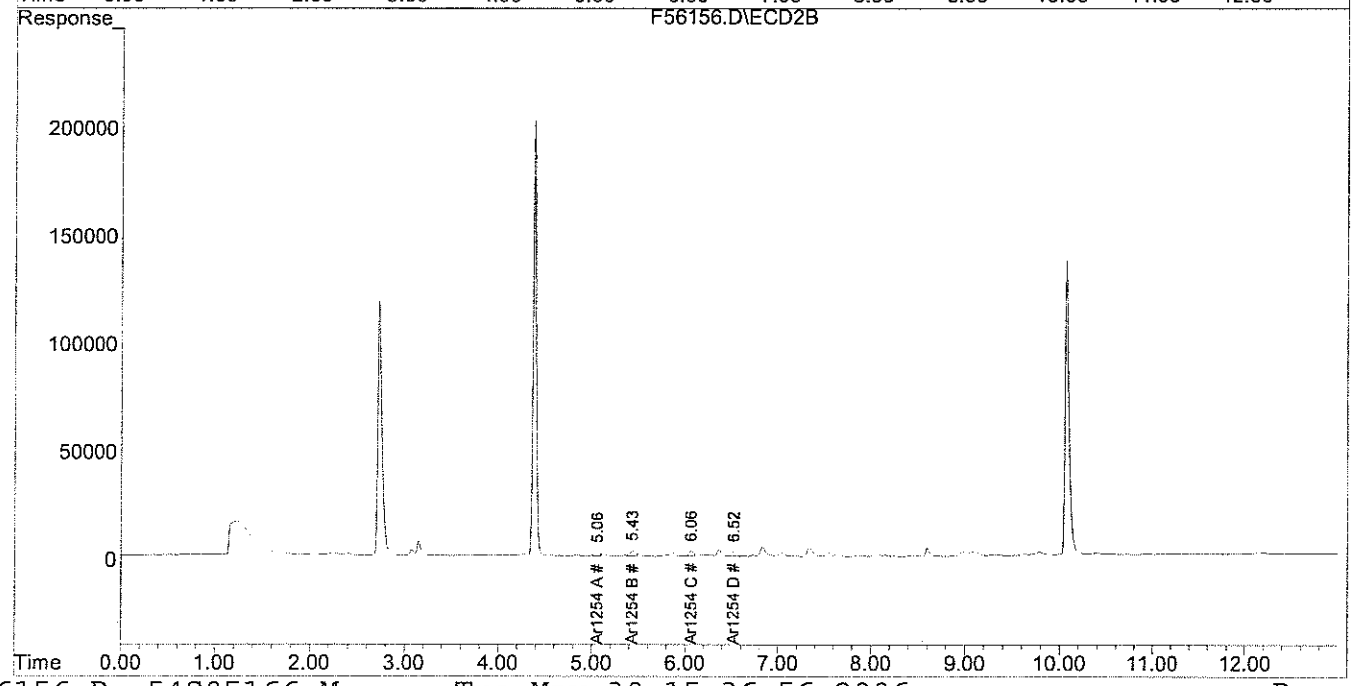
Quantitation Report (QT Reviewed)

Signal #1 : C:\HPCHEM\2\DATA\053006-F\F56156.D\ECD1A.CH Vial: 9
Signal #2 : C:\HPCHEM\2\DATA\053006-F\F56156.D\ECD2B.CH
Acq On : 30 May 2006 3:26 pm Operator:
Sample : 56637-6, A/C Inst : Instrumen
Misc : SOIL Multiplr: 1.00
IntFile Signal #1: EVENTS.E IntFile Signal #2: EVENTS2.E
Quant Time: May 30 15:36 2006 Quant Results File: 54S05166.RES

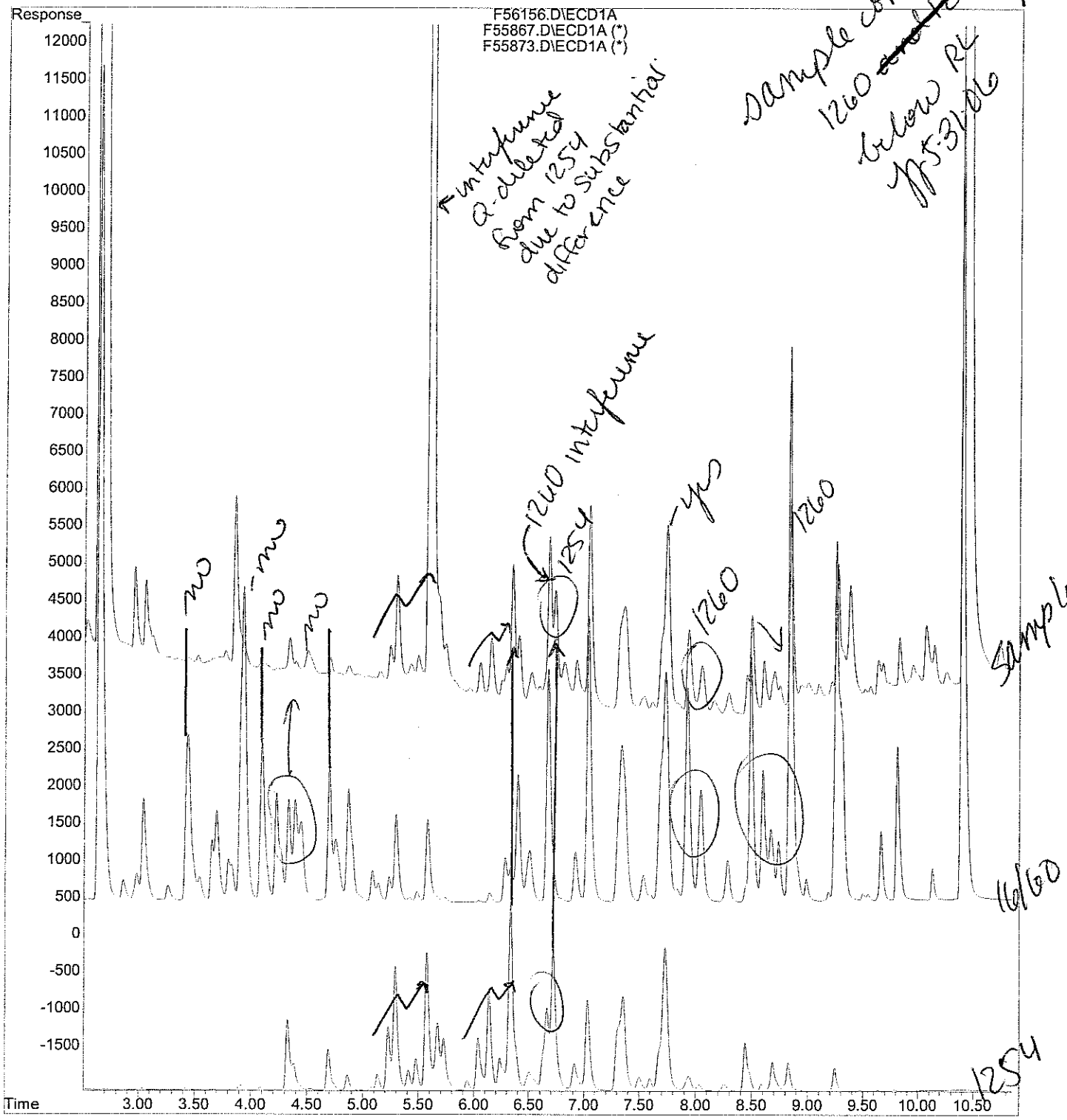
Quant Method : C:\HPCHEM\2\METHODS\54S05166.M (Chemstation Integrator)
Title : Aroclor 1254
Last Update : Fri May 26 08:03:09 2006
Response via : Single Level Calibration
DataAcq Meth : PSTQUICK.M

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film





File : C:\HPCHEM\2\DATA\053006-F\F56156.D
Operator :
Acquired : 30 May 2006 3:26 pm using AcqMethod PSTQUICK.M
Instrument : Instrumen
Sample Name: 56637-6, A/C
Misc Info : SOIL
Vial Number: 9



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Portland ME 04102

May 31, 2006

CLIENT SAMPLE ID

Project Name: GP Old Town Mill

Project Number: 218736.002

Field Sample ID: SB-35 (0-6")

SAMPLE DATA

Lab Sample ID: 56637-7
Matrix: Solid
Percent Solid: 96
Dilution Factor: 1.0
Collection Date: 05/25/06
Lab Receipt Date: 05/27/06
Extraction Date: 05/30/06
Analysis Date: 05/30/06

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit µg/kg	Results µg/kg
PCB-1016	15	U
PCB-1221	15	U
PCB-1232	15	U
PCB-1242	15	U
PCB-1248	15	U
PCB-1254	15	U
PCB-1260	15	U

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	87	%
Decachlorobiphenyl	82	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082.

Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3545.

COMMENTS: Results are expressed on a dry weight basis.

PCB Report

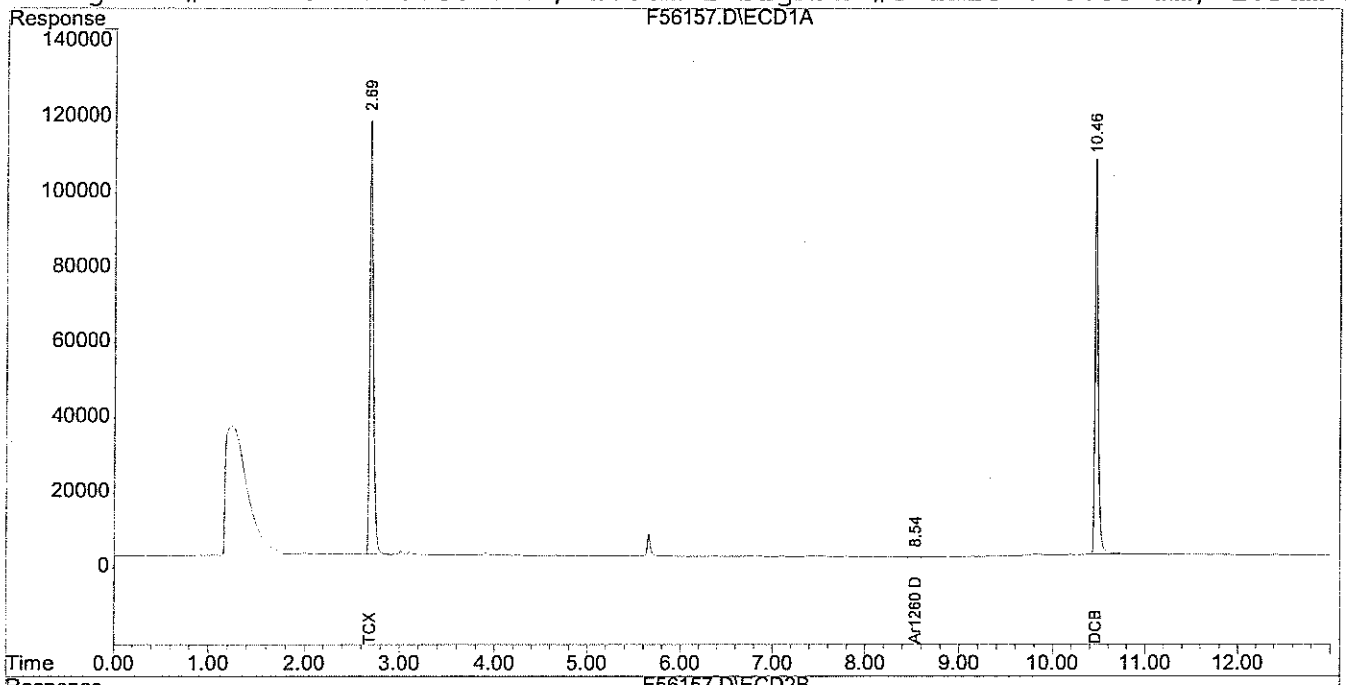
Authorized signature

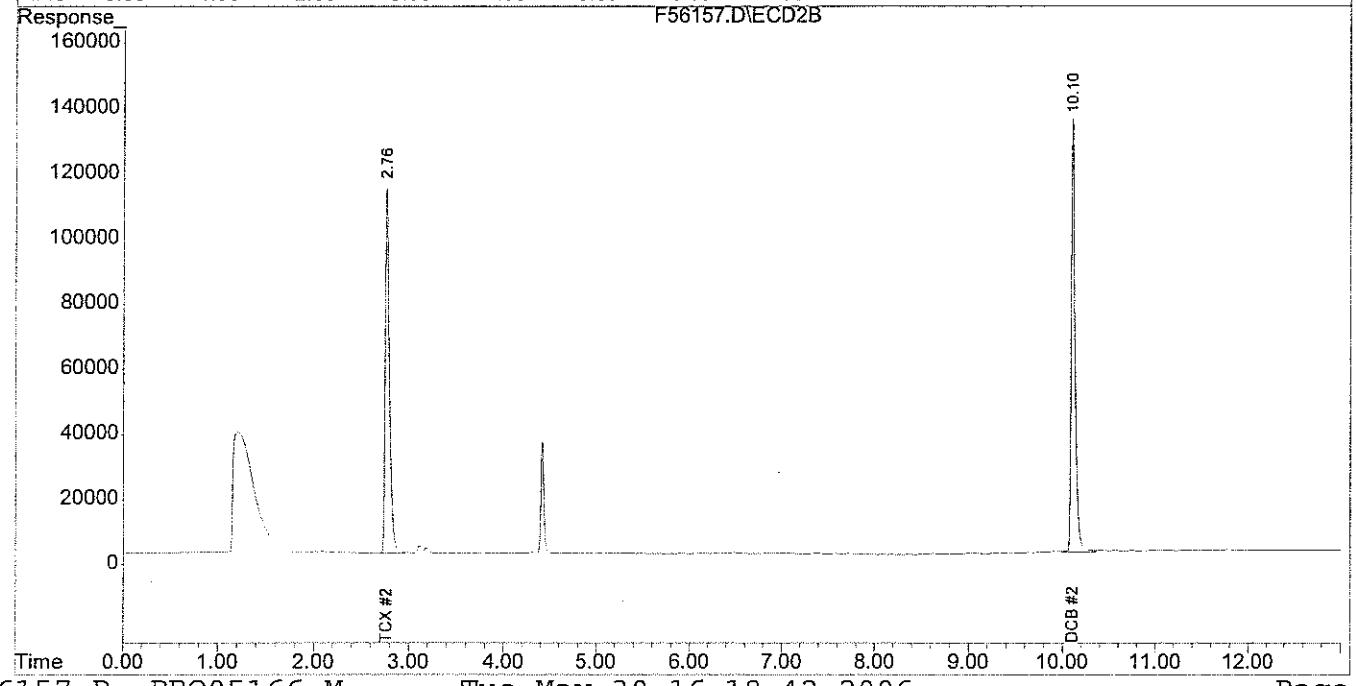
Quantitation Report (QT Reviewed)

Signal #1 : C:\HPCHEM\2\DATA\053006-F\F56157.D\ECD1A.CH Vial: 10
Signal #2 : C:\HPCHEM\2\DATA\053006-F\F56157.D\ECD2B.CH
Acq On : 30 May 2006 4:10 pm Operator:
Sample : 56637-7, A/C Inst : Instrumen
Misc : SOIL Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 30 16:19 2006 Quant Results File: PBQ05166.RES

Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 10:06:32 2006
Response via : Multiple Level Calibration
DataAcq Meth : PSTQUICK.M

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film





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Portland ME 04102

May 31, 2006

CLIENT SAMPLE ID

Project Name: GP Old Town Mill

Project Number: 218736.002

Field Sample ID: SB-5 (4-6')

SAMPLE DATA

Lab Sample ID: 56637-8
Matrix: Solid
Percent Solid: 92
Dilution Factor: 1.0
Collection Date: 05/25/06
Lab Receipt Date: 05/27/06
Extraction Date: 05/30/06
Analysis Date: 05/30/06

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit µg/kg	Results µg/kg
PCB-1016	15	U
PCB-1221	15	U
PCB-1232	15	U
PCB-1242	15	U
PCB-1248	15	U
PCB-1254	15	U
PCB-1260	15	U

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	95	%
Decachlorobiphenyl	85	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082.

Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3545.

COMMENTS: Results are expressed on a dry weight basis.

PCB Report

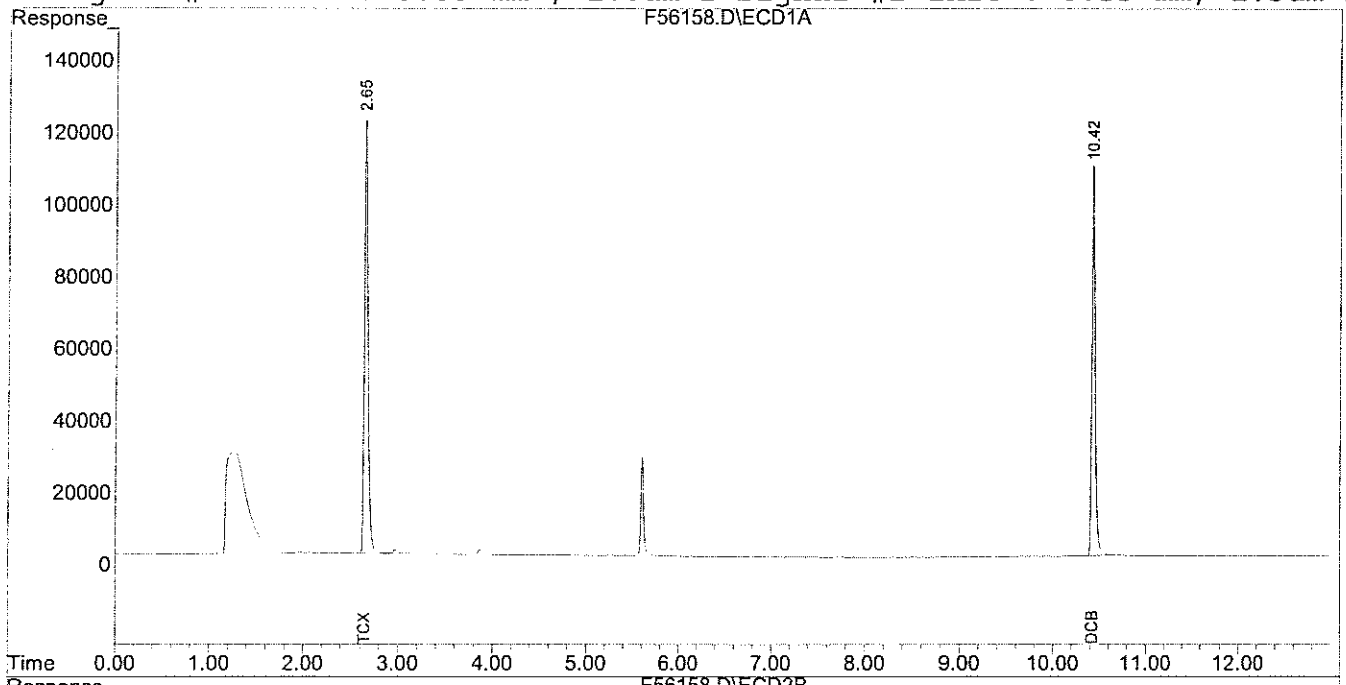
Authorized signature Melissa Nell

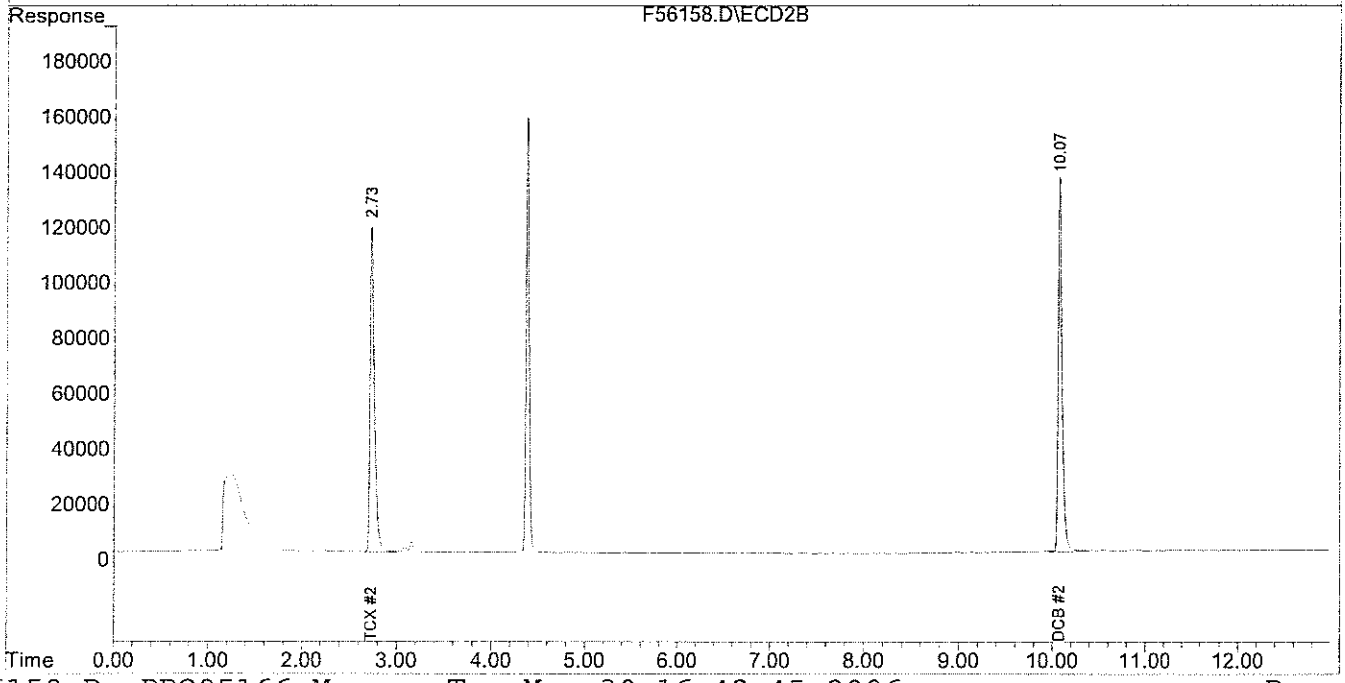
Quantitation Report (QT Reviewed)

Signal #1 : C:\HPCHEM\2\DATA\053006-F\F56158.D\ECD1A.CH Vial: 11
Signal #2 : C:\HPCHEM\2\DATA\053006-F\F56158.D\ECD2B.CH
Acq On : 30 May 2006 4:24 pm Operator:
Sample : 56637-8, A/C Inst : Instrumen
Misc : SOIL Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 30 16:42 2006 Quant Results File: PBQ05166.RES

Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 10:06:32 2006
Response via : Multiple Level Calibration
DataAcq Meth : PSTQUICK.M

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film





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May 31, 2006

CLIENT SAMPLE ID

Project Name: GP Old Town Mill

Project Number: 218736.002

Field Sample ID: SB-21 (0-4')

SAMPLE DATA

Lab Sample ID: 56637-9
Matrix: Solid
Percent Solid: 97
Dilution Factor: 1.0
Collection Date: 05/25/06
Lab Receipt Date: 05/27/06
Extraction Date: 05/30/06
Analysis Date: 05/30/06

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit µg/kg	Results µg/kg
PCB-1016	15	U
PCB-1221	15	U
PCB-1232	15	U
PCB-1242	15	U
PCB-1248	15	U
PCB-1254	15	U
PCB-1260	15	U

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	94	%
Decachlorobiphenyl	89	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082.

Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3545.

COMMENTS: Results are expressed on a dry weight basis.

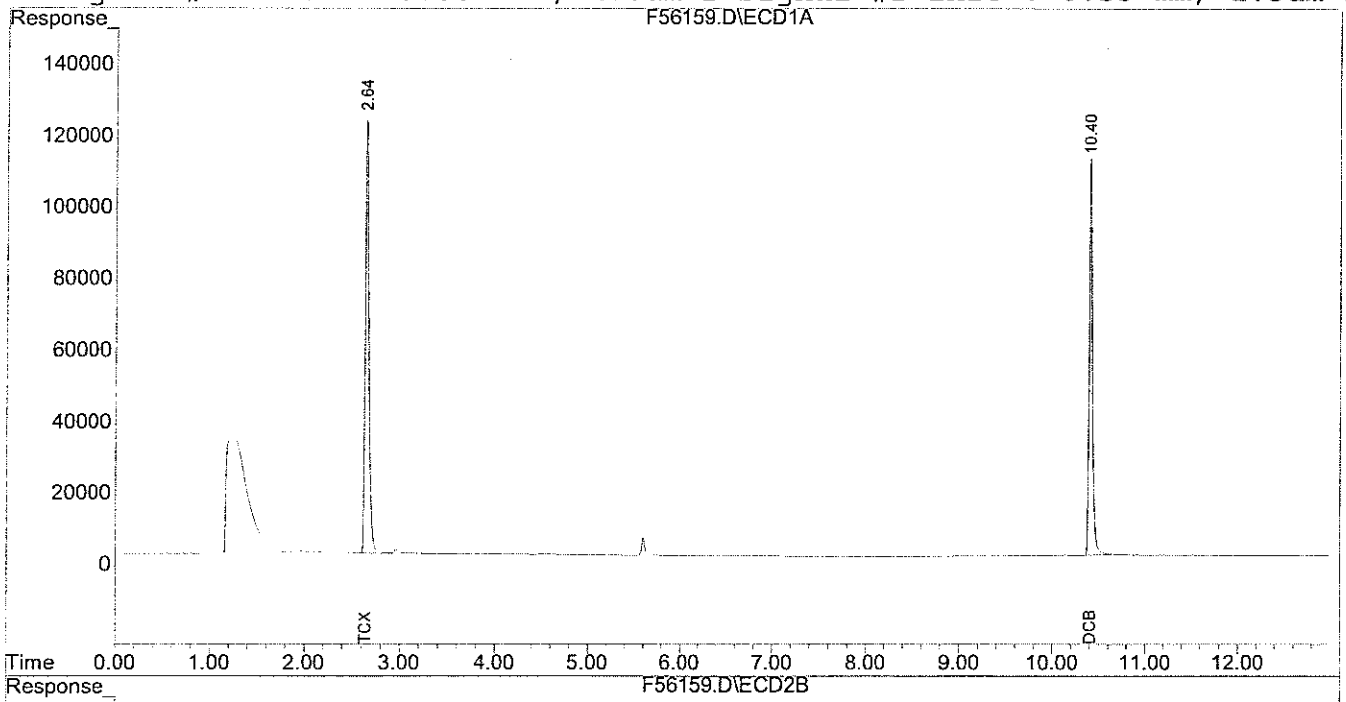
PCB Report

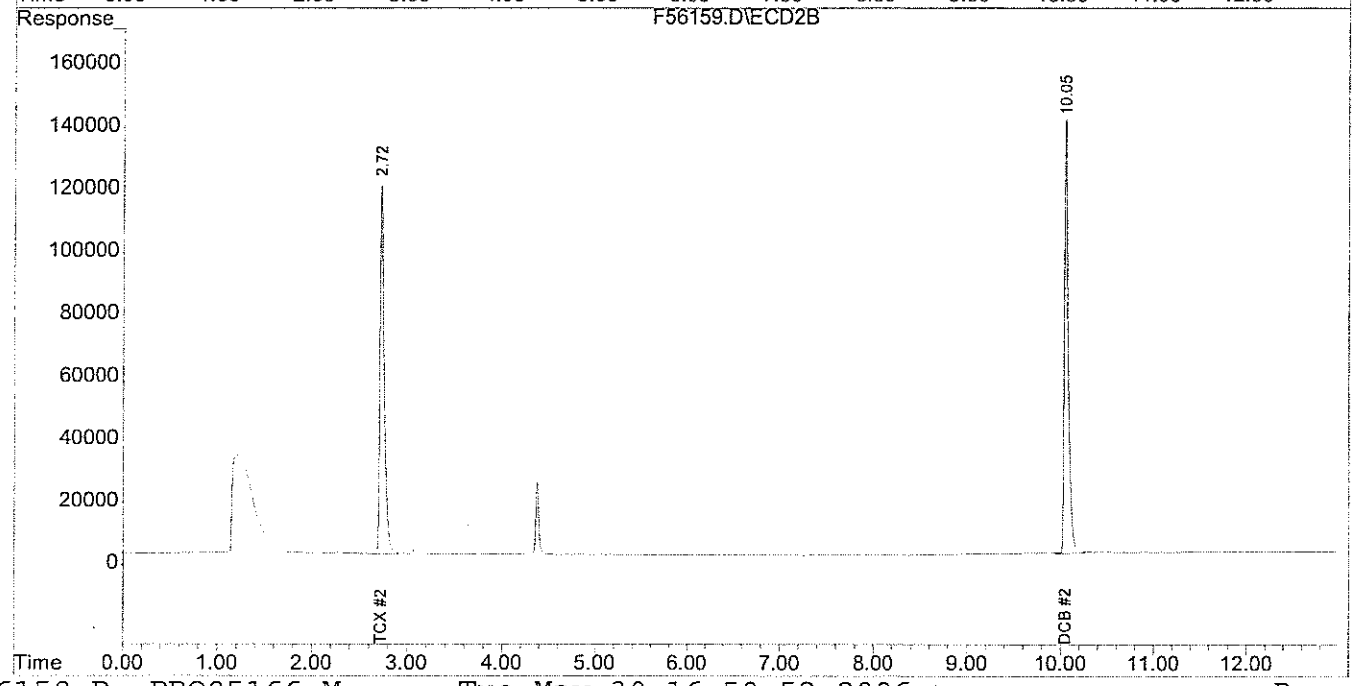
Authorized signature

Signal #1 : C:\HPCHEM\2\DATA\053006-F\F56159.D\ECD1A.CH Vial: 12
Signal #2 : C:\HPCHEM\2\DATA\053006-F\F56159.D\ECD2B.CH
Acq On : 30 May 2006 4:39 pm Operator:
Sample : 56637-9, A/C Inst : Instrumen
Misc : SOIL Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 30 16:50 2006 Quant Results File: PBQ05166.RES

Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 10:06:32 2006
Response via : Multiple Level Calibration
DataAcq Meth : PSTQUICK.M

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film





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May 31, 2006

CLIENT SAMPLE ID

Project Name: GP Old Town Mill

Project Number: 218736.002

Field Sample ID: SB-5 (0-4')

SAMPLE DATA

Lab Sample ID: 56637-10
Matrix: Solid
Percent Solid: 94
Dilution Factor: 1.0
Collection Date: 05/25/06
Lab Receipt Date: 05/27/06
Extraction Date: 05/30/06
Analysis Date: 05/30/06

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit µg/kg	Results µg/kg
PCB-1016	15	U
PCB-1221	15	U
PCB-1232	15	U
PCB-1242	15	U
PCB-1248	15	U
PCB-1254	15	U
PCB-1260	15	U

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	93	%
Decachlorobiphenyl	87	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082.

Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3545.

COMMENTS: Results are expressed on a dry weight basis.

PCB Report

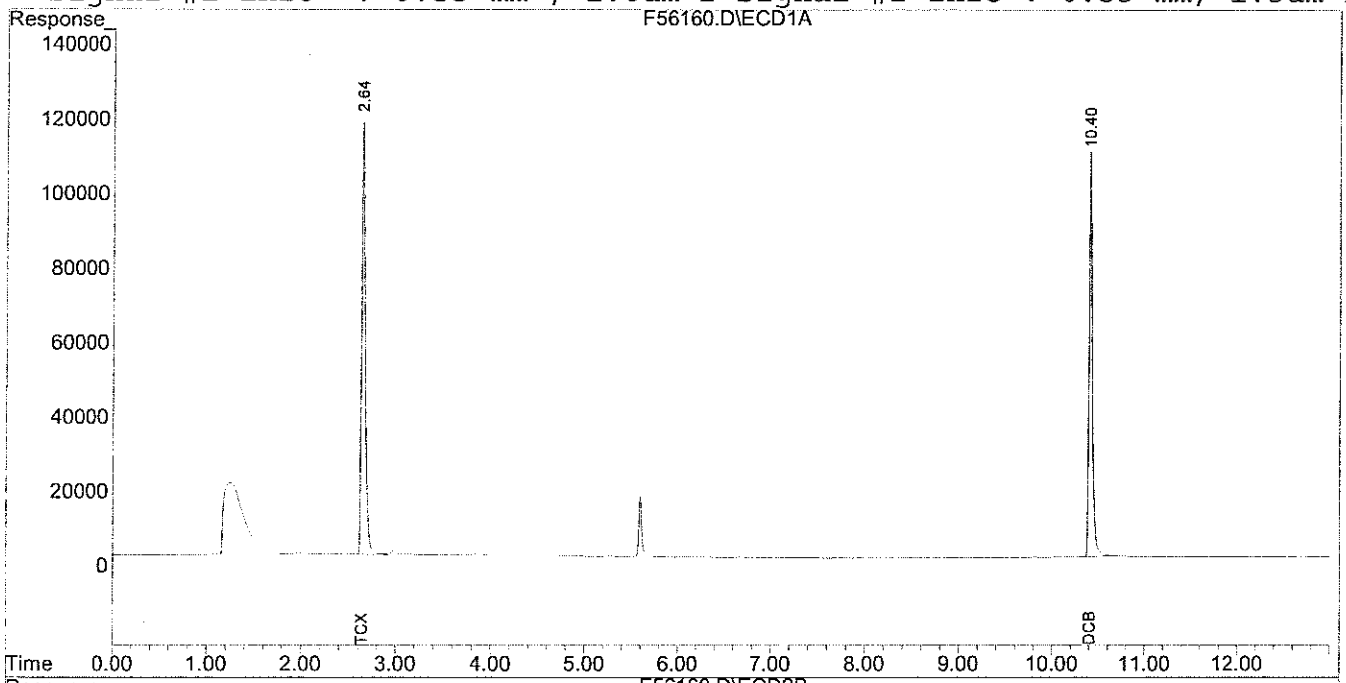
Authorized signature Melissa Sull

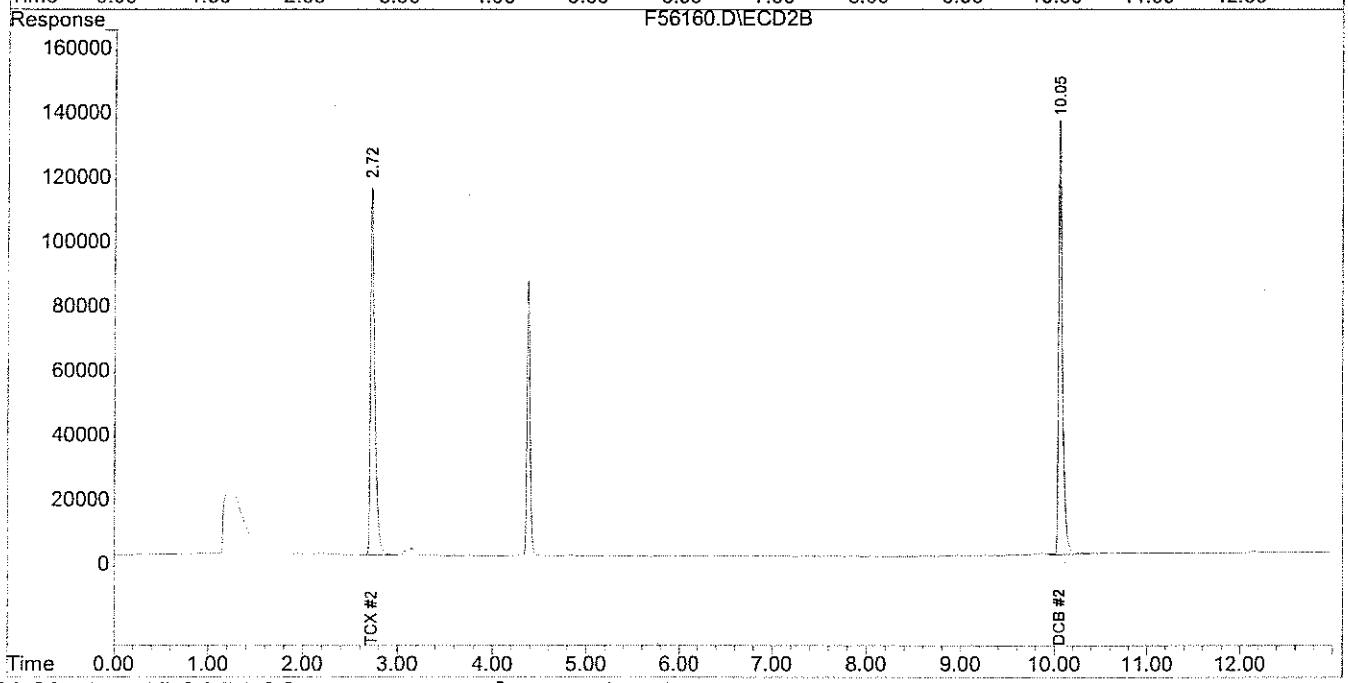
Quantitation Report (QT Reviewed)

Signal #1 : C:\HPCHEM\2\DATA\053006-F\F56160.D\ECD1A.CH Vial: 13
Signal #2 : C:\HPCHEM\2\DATA\053006-F\F56160.D\ECD2B.CH
Acq On : 30 May 2006 4:53 pm Operator:
Sample : 56637-10, A/C Inst : Instrumen
Misc : SOIL Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 31 8:03 2006 Quant Results File: PBQ05166.RES

Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 10:06:32 2006
Response via : Multiple Level Calibration
DataAcq Meth : PSTQUICK.M

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film





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May 31, 2006

CLIENT SAMPLE ID

Project Name: GP Old Town Mill

Project Number: 218736.002

Field Sample ID: SB-23 (0-4')

SAMPLE DATA

Lab Sample ID: 56637-12
Matrix: Solid
Percent Solid: 90
Dilution Factor: 1.1
Collection Date: 05/25/06
Lab Receipt Date: 05/27/06
Extraction Date: 05/30/06
Analysis Date: 05/30/06

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit µg/kg	Results µg/kg
PCB-1016	17	U
PCB-1221	17	U
PCB-1232	17	U
PCB-1242	17	U
PCB-1248	17	U
PCB-1254	17	U
PCB-1260	17	U

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	89	%
Decachlorobiphenyl	84	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082.

Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3545.

COMMENTS: Results are expressed on a dry weight basis.

PCB Report

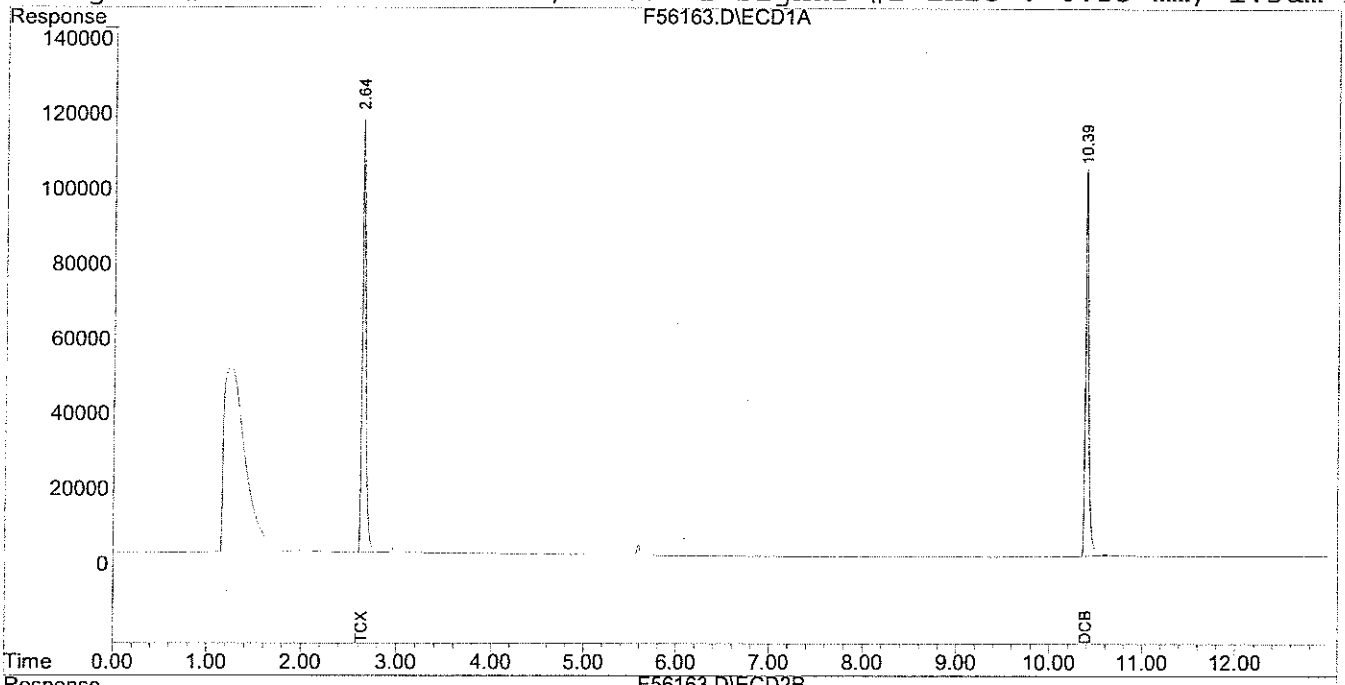
Authorized signature

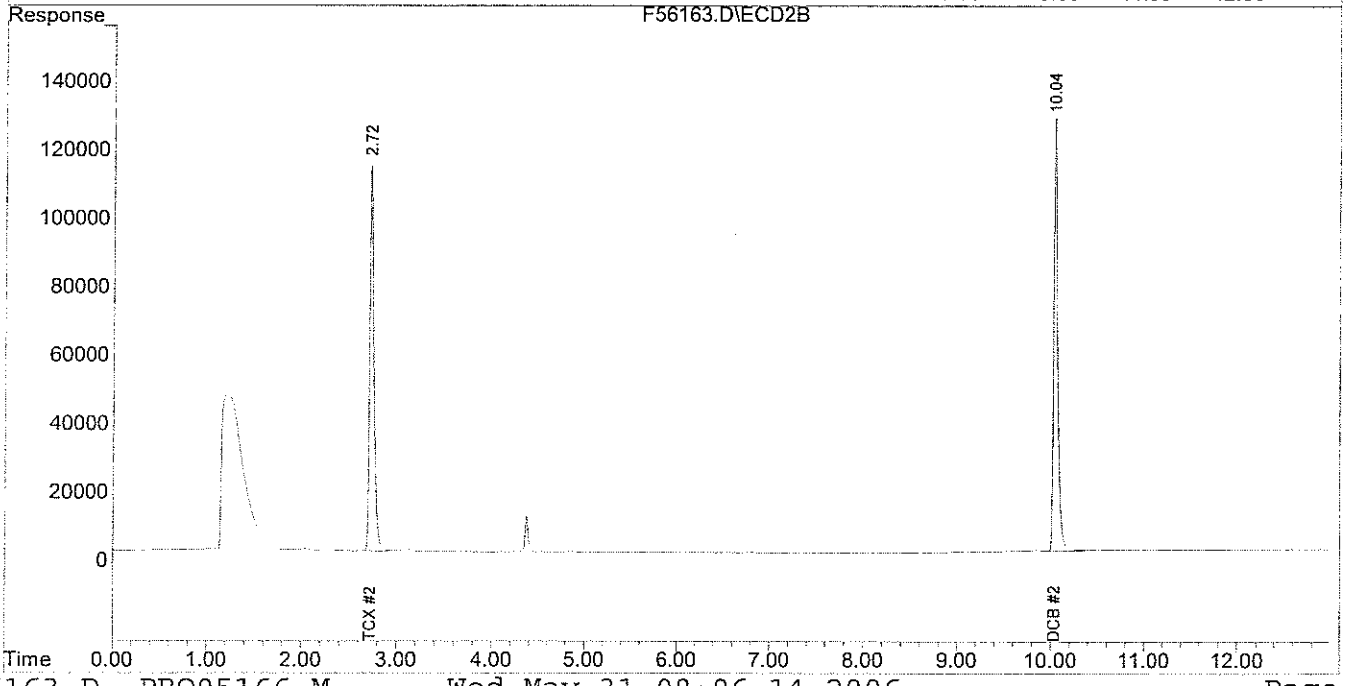
Quantitation Report (QT Reviewed)

Signal #1 : C:\HPCHEM\2\DATA\053006-F\F56163.D\ECD1A.CH Vial: 16
Signal #2 : C:\HPCHEM\2\DATA\053006-F\F56163.D\ECD2B.CH
Acq On : 30 May 2006 5:52 pm Operator:
Sample : 56637-12, A/C Inst : Instrumen
Misc : SOIL Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 31 8:06 2006 Quant Results File: PBQ05166.RES

Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 10:06:32 2006
Response via : Multiple Level Calibration
DataAcq Meth : PSTQUICK.M

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film





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May 31, 2006

CLIENT SAMPLE ID

Project Name: GP Old Town Mill

Project Number: 218736.002

Field Sample ID: SB-24 (0-4')

SAMPLE DATA

Lab Sample ID: 56637-13
Matrix: Solid
Percent Solid: 76
Dilution Factor: 1.2
Collection Date: 05/25/06
Lab Receipt Date: 05/27/06
Extraction Date: 05/30/06
Analysis Date: 05/30/06

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit µg/kg	Results µg/kg
PCB-1016	18	U
PCB-1221	18	U
PCB-1232	18	U
PCB-1242	18	U
PCB-1248	18	U
PCB-1254	18	U
PCB-1260	18	U

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	65	%
Decachlorobiphenyl	72	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082.

COMMENTS: Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3545. Results are expressed on a dry weight basis.

PCB Report

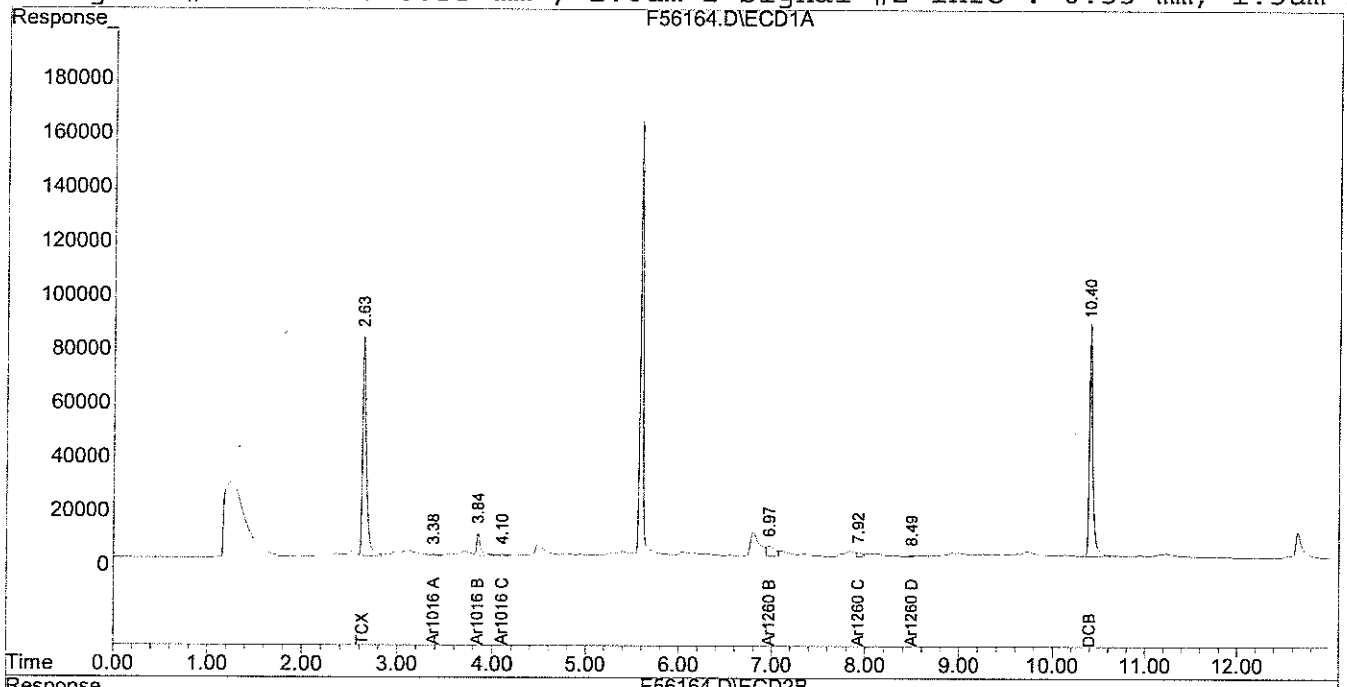
Authorized signature

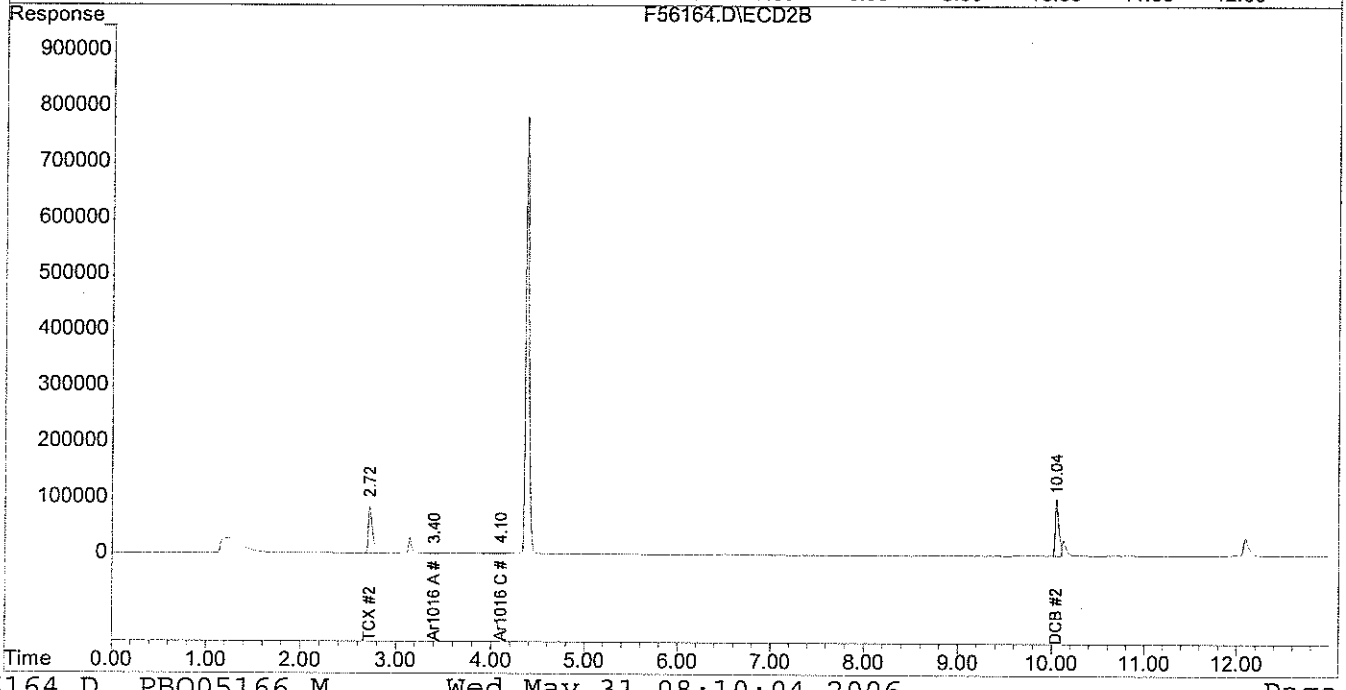
Quantitation Report (QT Reviewed)

Signal #1 : C:\HPCHEM\2\DATA\053006-F\F56164.D\ECD1A.CH Vial: 17
Signal #2 : C:\HPCHEM\2\DATA\053006-F\F56164.D\ECD2B.CH
Acq On : 30 May 2006 6:07 pm Operator:
Sample : 56637-13, A/C Inst : Instrumen
Misc : SOIL Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 31 8:07 2006 Quant Results File: PBQ05166.RES

Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 10:06:32 2006
Response via : Multiple Level Calibration
DataAcq Meth : PSTQUICK.M

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film





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Portland ME 04102

May 31, 2006

CLIENT SAMPLE ID

Project Name: GP Old Town Mill

Project Number: 218736.002

Field Sample ID: SB-41 (0-4)

SAMPLE DATA

Lab Sample ID: 56637-19
Matrix: Solid
Percent Solid: 91
Dilution Factor: 22
Collection Date: 05/26/06
Lab Receipt Date: 05/27/06
Extraction Date: 05/30/06
Analysis Date: 05/31/06

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit µg/kg	Results µg/kg
PCB-1016	330	U
PCB-1221	330	U
PCB-1232	330	U
PCB-1242	330	U
PCB-1248	330	U
PCB-1254	330	U
PCB-1260	330	**4050**

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	*	%
Decachlorobiphenyl	*	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082.

COMMENTS: Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3545. Results are expressed on a dry weight basis. * The surrogates were diluted out.

PCB Report

Authorized signature

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PCB CONFIRMATION RESULTS

SDG: 56637

Project No: 218736.002 Site ID: Location: ME

Instrument ID: F GC Columns: RTX-CLPesticidesI ID: 0.32(mm)

RTX-CLPesticidesII

Sample ID: 56637-19

COMPOUND	RESPONSE #1 ug/kg	RESPONSE #2 ug/kg	DIFFERENCE %	#	% DIFFERENCE Acceptance Limit
PCB 1260	4050	4370	8		40

Column to be used to flag % Difference with an asterisk

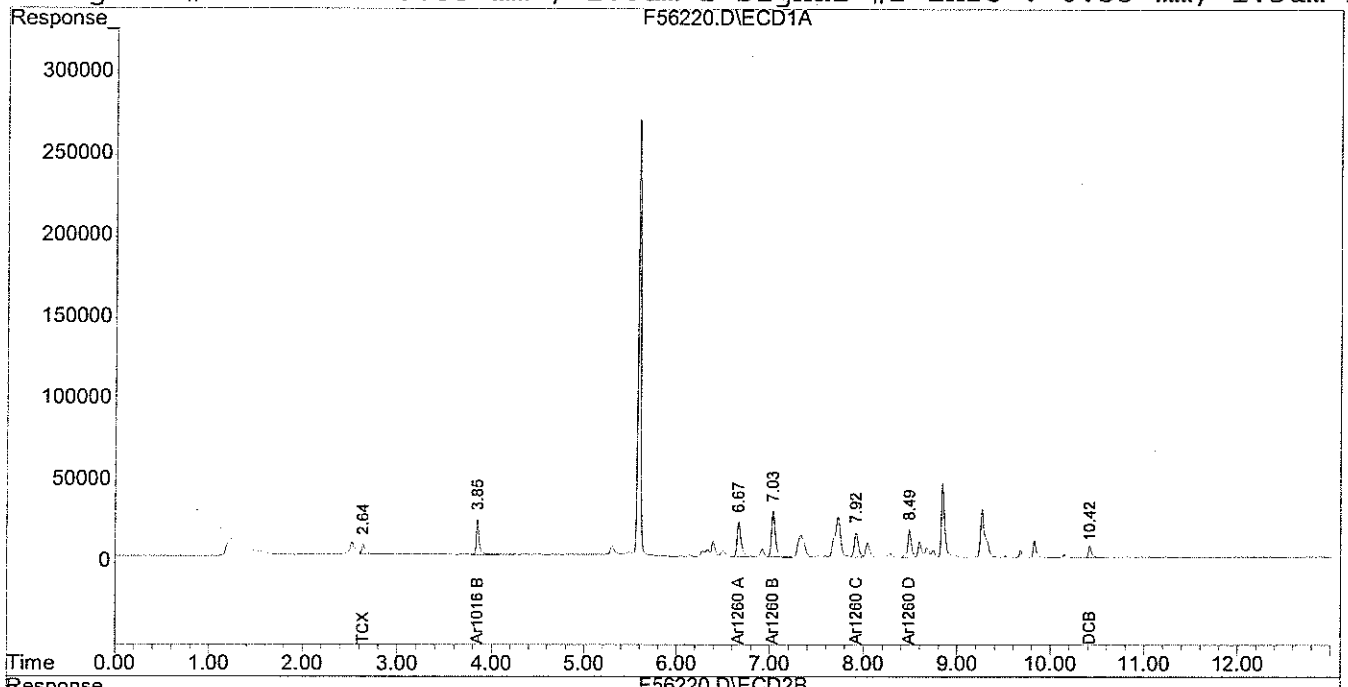
* Values outside of QC limits

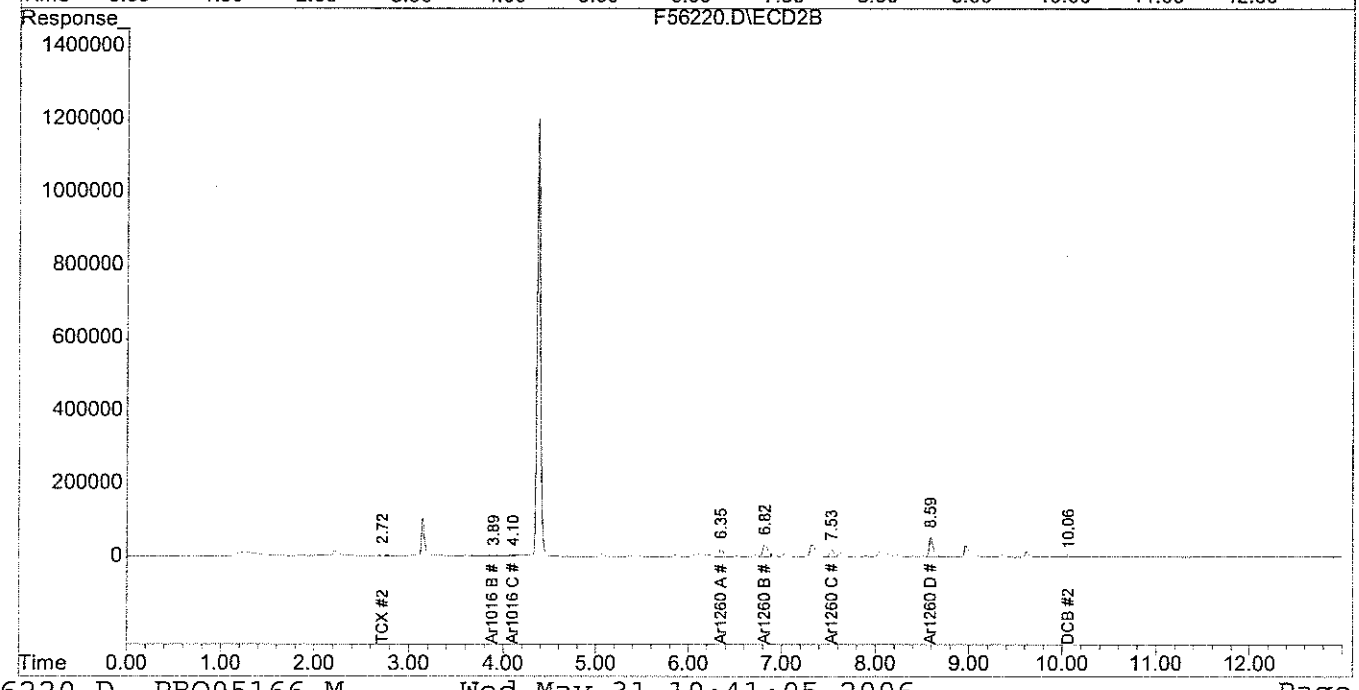
Quantitation Report (QT Reviewed)

Signal #1 : C:\HPCHEM\2\DATA\053106-F\F56220.D\ECD1A.CH Vial: 3
Signal #2 : C:\HPCHEM\2\DATA\053106-F\F56220.D\ECD2B.CH
Acq On : 31 May 2006 10:12 am Operator:
Sample : 56637-19, 1:20, A/C Inst : Instrumen
Misc : SOIL Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 31 10:40 2006 Quant Results File: PBQ05166.RES

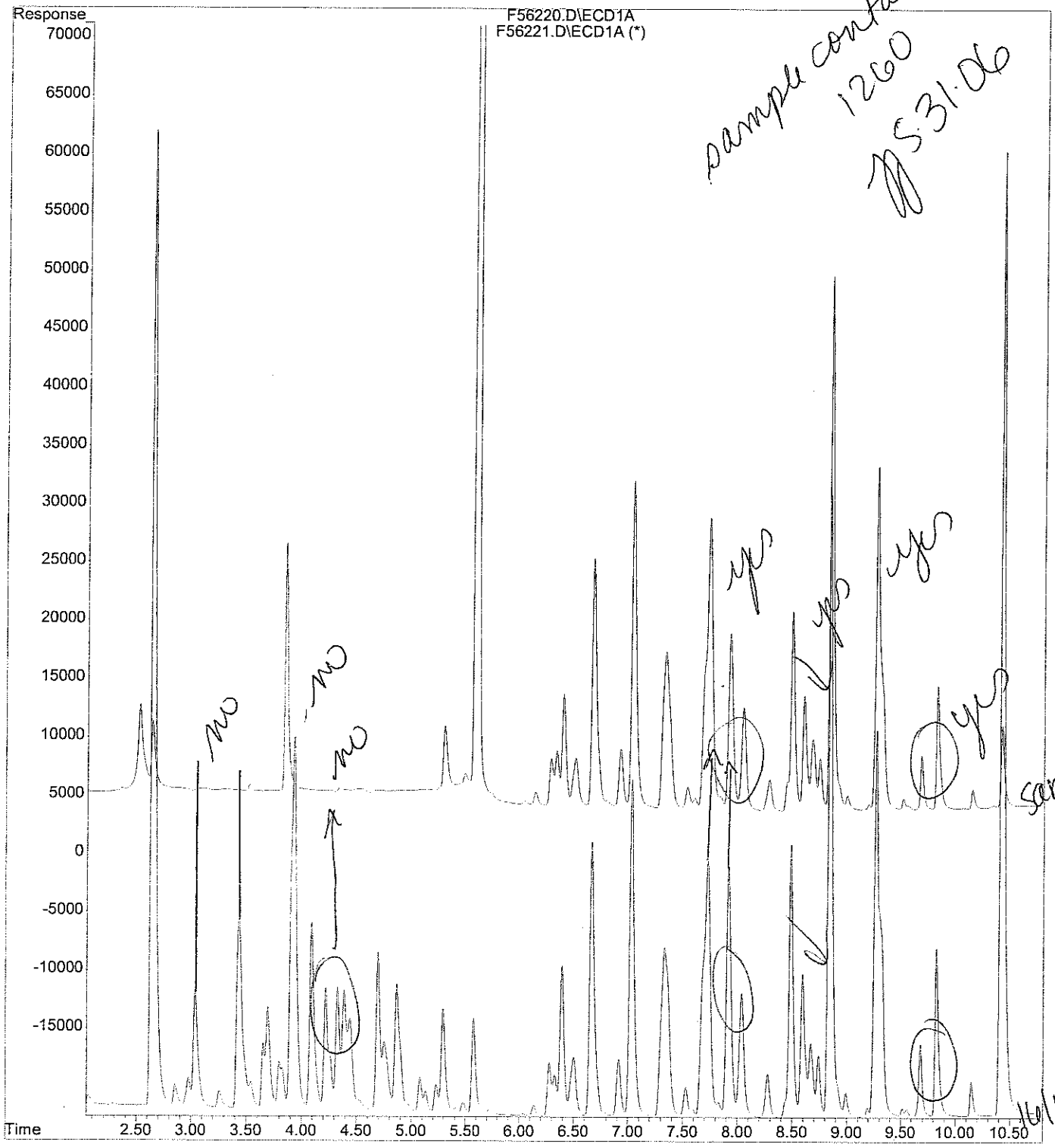
Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 10:06:32 2006
Response via : Multiple Level Calibration
DataAcq Meth : PSTQUICK.M

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film





File : C:\HPCHEM\2\DATA\053106-F\F56220.D
Operator :
Acquired : 31 May 2006 10:12 am using AcqMethod PSTQUICK.M
Instrument : Instrumen
Sample Name: 56637-19, 1:20, A/C
Misc Info : SOIL
Vial Number: 3



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PCB
QC FORMS

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PCB SOLID
SYSTEM MONITORING COMPOUND RECOVERY

Project #: 218736.002 Quote #: SDG#: 56637

Site ID: GP Old Town Mill Location: ME

Instrument ID: F
GC Column: RTX-CLPesticides I ID: 0.32(mm)
RTX-CLPesticides II

	SAMPLE ID	COLUMN #1				COLUMN #2				TOTAL	
		SMC1	#	SMC2	#	SMC1	#	SMC2	#	OUT	#
01	P4016 16/60 0.01	NA		NA		NA		NA		0	
02	P4017 16/60 0.02	NA		NA		NA		NA		0	
03	P4018 16/60 0.05	NA		NA		NA		NA		0	
04	P4019 16/60 0.1	NA		NA		NA		NA		0	
05	P4020 16/60 0.25	116		109		116		121		0	
06	P4021 16/60 0.5	112		104		110		108		0	
07	P4022 16/60 1.0	NA		NA		NA		NA		0	
08	P4024 16/60 0.5 CC	104		104		105		107		0	
09	P3883 AR 1221	NA		NA		NA		NA		0	
10	P3884 AR 1232	NA		NA		NA		NA		0	
11	P3885 AR 1242	NA		NA		NA		NA		0	
12	P3886 AR 1248	NA		NA		NA		NA		0	
13	P3994 AR 1254	NA		NA		NA		NA		0	
14	P3887 AR 1262	NA		NA		NA		NA		0	
15	P3888 AR 1268	NA		NA		NA		NA		0	
16	P4057 AR 1254 0.02	NA		NA		NA		NA		0	
17	P4056 AR 1254 0.05	NA		NA		NA		NA		0	
18	P4055 AR 1254 0.1	NA		NA		NA		NA		0	
19	P4054 AR 1254 0.25	NA		NA		NA		NA		0	
20	P4053 AR 1254 0.5	NA		NA		NA		NA		0	
21	P4052 AR 1254 1.0	NA		NA		NA		NA		0	
22	P4010 AR 1254 0.5 C	102		103		105		108		0	
23	ZZZZ	81		17	*	87		14	*	2	
24	ZZZZ	85		21	*	95		11	*	2	
25	ZZZZ	87		15	*	95		16	*	2	
26	ZZZZ	86		14	*	94		15	*	2	
27	ZZZZ	77		81		80		81		0	
28	ZZZZ	62		33		87		33		0	
29	ZZZZ	83		10	*	96		9	*	2	
30	P4021 16/60 0.5	99		103		102		109		0	
31	P4053 AR 1254 0.5	NA		NA		NA		NA		0	
32											
33											
34											
35											
36											
37											
38											
39											
40											
41											
42											

QC LIMITS
SMC1 = TCX 46-122
SMC2 = DCB 26-86

*=Values outside of protocol required QC limits
D=System Monitoring Compound diluted out

analytics environmental laboratory LLC

PCB SOLID
SYSTEM MONITORING COMPOUND RECOVERY

Project #: 218736.002 Quote #: SDG#: 56637

Site ID: GP Old Town Mill Location: ME

Instrument ID: F
GC Column: RTX-CLPesticides I ID: 0.32(mm)
RTX-CLPesticides II

		COLUMN #1			COLUMN #2			TOTAL	
	SAMPLE ID	SMC1	#	SMC2	#	SMC1	#	SMC2	#	OUT	#
01	P4021 16/60 0.5	93		89		96		92		0	
02	B05306PAS, A/C	64		44		83		83		0	
03	L05306PASB, A/C	95		76		101		96		0	
04	LD05306PASB, A/C	92		81		97		93		0	
05	56637-4, A/C	92		81		97		91		0	
06	56637-6, A/C	93		86		98		94		0	
07	56637-7, A/C	87		82		93		93		0	
08	56637-8, A/C	95		85		99		94		0	
09	56637-9, A/C	94		89		98		96		0	
10	56637-10, A/C	93		87		97		94		0	
11	56637-10, MS, A/C	96		90		99		95		0	
12	56637-10, MSD, A/C	95		89		99		94		0	
13	56637-12, A/C	89		84		93		89		0	
14	56637-13, A/C	65		72		69		62		0	
15	ZZZZZ	91		73		95		67		0	
16	P4021 16/60 0.5	88		85		93		87		0	
17											
18											
19											
20											
21											
22											
23											
24											
25											
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37											
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39											
40											
41											
42											

QC LIMITS
SMC1 = TCX 46-123
SMC2 = DCB 38-123

*=Values outside of protocol required QC limits
D=System Monitoring Compound diluted out

analytics environmental laboratory LLC

PCB SOLID
SYSTEM MONITORING COMPOUND RECOVERY

Project #: 218736.002 Quote #: SDG#: 56637

Site ID: GP Old Town Mill Location: ME

Instrument ID: F
GC Column: RTX-CLPesticides I ID: 0.32(mm)
RTX-CLPesticides II

		COLUMN #1			COLUMN #2			TOTAL	
	SAMPLE ID	SMC1	#	SMC2	#	SMC1	#	SMC2	#	OUT	#
01	P4020 16/60 0.25	98		106		100		102		0	
02	B05306PAS,RR A/C	88		92		88		93		0	
03	56637-19, 1:20, A/C	D		D		D		D		0	
04	P4021 16/60 0.5	99		100		100		101		0	
05											
06											
07											
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42											

QC LIMITS
SMC1 = TCX 46-123
SMC2 = DCB 38-123
*=Values outside of protocol required QC limits
D=System Monitoring Compound diluted out

analytics environmental laboratory LLC

PCB SOLID
LABORATORY CONTROL SPIKE/ DUPLICATE
PERCENT RECOVERY

Project #: 218736.002 Quote #: SDG#: 56637

Site ID: GP Old Town Mill Location: ME

Instrument ID: F

GC Column: RTX-CLPesticides I ID: 0.32(mm)

RTX-CLPesticides II

Laboratory Control Spike/Duplicate Sample ID: L05306PASB A/C/ LD05306PASB A/C

COMPOUND	SPIKE ADDED (ug/kg)	SAMPLE CONCENTRATION (ug/kg)	LCS CONCENTRATION (ug/kg)	LCS RECOVERY % #	QC LIMITS % REC
PCB 1016	267	U	243	91	67-144
PCB 1260	267	U	219	82	58-155
PCB 1016 #2	267	U	265	99	69-140
PCB 1260 #2	267	U	278	104	67-149

COMPOUND	SPIKE ADDED (ug/kg)	LCSD CONCENTRATION (ug/kg)	LCSD % REC #	RPD #	QC LIMITS RPD	QC LIMITS % REC.
PCB 1016	267	245	92	1	30	67-144
PCB 1260	267	224	84	2	30	58-155
PCB 1016 #2	267	262	98	1	30	69-140
PCB 1260 #2	267	271	102	2	30	67-149

Column to be used to flag recovery and RPD values with an asterisk

* Values outside of QC limits
RPD: 0 out of 4 outside limits
Spike Recovery: 0 out of 8 outside limits

Comments: _____

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PCB SOLID
MATRIX SPIKE/ DUPLICATE
PERCENT RECOVERY

Project #: 216736.002 Quote #: SDG#: 56637

Site ID: GP Old Town Mill Location: ME

Instrument ID: F

GC Column: RTX-CL Pesticides I ID: 0.32(mm)

RTX-CL Pesticides II

Matrix Spike/Duplicate Sample ID: 56637-10

COMPOUND	SPIKE ADDED (ug/kg)	SAMPLE CONCENTRATION (ug/kg)	MS CONCENTRATION (ug/kg)	MS RECOVERY %	#	QC LIMITS % REC
PCB 1016	281	U	272	97		67-144
PCB 1260	281	U	265	95		58-155
PCB 1016 #2	281	U	282	101		69-140
PCB 1260 #2	281	U	289	103		67-149

COMPOUND	SPIKE ADDED (ug/kg)	MSD CONCENTRATION (ug/kg)	MSD % REC	#	RPD	#	QC LIMITS RPD	QC LIMITS % REC.
PCB 1016	276	247	90		10		30	67-144
PCB 1260	276	239	87		10		30	58-155
PCB 1016 #2	276	258	94		9		30	69-140
PCB 1260 #2	276	264	96		9		30	67-149

Column to be used to flag recovery and RPD values with an asterisk

* Values outside of QC limits
RPD: 0 out of 4 outside limits
Spike Recovery: 0 out of 8 outside limits

Comments: _____

analytics environmental laboratory LLC

PEST/PCB/HERB
Instrument F
Sequence Log Book

051606 -F

Load: JJ
Unload: JJ

Data File	Vial #	Sample	Method	Comments
F55857	9	P4045 Pest50		✓
58	10	P4046 Pest100		✓
59	11	P4047 Pest150		✓
60	12	P3965 Pest 0.5CC		✓
61	13	P4016 16/60 0.01	PBQ05166	✓
62	14	P4017 16/60 0.02		✓
63	15	P4018 16/60 0.05		✓
64	16	P4019 16/60 0.1		✓
65	17	P4020 16/60 0.25		✓
66	18	P4021 16/60 0.5		✓
67	19	P4022 16/60 1.0		✓
68	20	P4024 16/60 0.5CC		✓
69	21	P3883 AR1221		for overlay, single point
70	22	P3884 AR1232		
71	23	P3885 AR 1242		
72	24	P3886 AR 1248		
73	25	P3994 AR 1254		
74	26	P3887 AR 1262		
75	27	P3888 AR 1268		
76	28	P4057 AR1254 0.02	JJ 5.18.06 ~~P~~54Q05166	✓
77	29	P4056 AR1254 0.05		✓
78	30	P4055 AR1254 0.1		✓
79	31	P4054 AR1254 0.25		✓
80	32	P4053 AR1254 0.5		✓
81	33	P4052 AR1254 1.0		✓
82	34	P4010 AR1254 0.5		✓
83	35	B05106PW		✓
84	36	L05106PWT		✓ methoxy ↑#1, ok#2, endrin
85	37	L05106PWB		✓ ~~AA~~ ketone ↑(marg.)#1, JJ 5.17.06 ↑#2

Standard Dilutions: 1:2- 500 uL sample +500 uL hexane
1:5- 200 uL sample +800 uL hexane
Hexane Lot : _____ 1:10- 100 uL sample + 900 uL hexane
1:20- 50 uL sample + 950 uL hexane
MTBE Lot: _____ 1:100- 10 uLsample +990 uL hexane
1:1000- 1 uLsample +1000 uL hexane

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PEST/PCB/HERB
Instrument F
Sequence Log Book

053006-F

Load: JG
Unload: JG

Data File	Vial #	Sample	Method	Comments
F56147	100	Hexane		✓
48	1	P4020 16/60 0.25	PBQ05166	✓
49	2	B05236PAS, A/C		✓
50	3	56599-1, 1:10 ↓		handed in w/sample
51	4	P4021 16/60 0.5		✓
52	5	B05306PAS A/C		✓
53	6	L05306PASB		✓
54	7	L05306PASB		✓
55	8	56637-4		✓
56	9	-6	54505166	✓ 1254 ↓ hit
57	10	-~~8~~ 7 A/C JG 5.30.06		✓
58	11	-8		✓
59	12	-9		✓
60	13	-10		✓
61	14	-10, MS A/C		✓
62	15	-10, MSD		✓
63	16	-12		✓
64	17	-13		✓
65	18	-19		RR@1:20
66	4	P4021 16/60 0.5		✓
67-27	100	Hexane		

JG 5.31.06

Standard Dilutions: 1:2- 500 uL sample +500 uL hexane
1:5- 200 uL sample +800 uL hexane
Hexane Lot : 46060 1:10- 100 uL sample + 900 uL hexane
1:20- 50 uL sample + 950 uL hexane
MTBE Lot: _____ 1:100- 10 uLsample +990 uL hexane
1:1000- 1 uLsample +1000 uL hexane

analytics environmental laboratory LLC

PEST/PCB/HERB
Instrument F
Sequence Log Book

053106-F

Load: yyy
Unload: yyy

Data File	Vial #	Sample	Method	Comments
F.56218	1	P4D2O 16/60 0.25	PBQ051606	✓
19	2	B053O6PAS, A/C RR		✓
20	3	56637-19, 1:20, A/C		✓
21	4	P4021 16/60 0.5		✓
22	100	Hexane		
23	99	P3862 PEM	PTQ051606	✓
24	1	P4044 Pest 30	PBQ051606	✓ DCB ↑ #2
25	2	P3981 C/T 0.25/0.5	CTQ05236	✓
26	3	P4020 16/60 0.25		✓
27	4	B053160PW		✓
28	5	L053160PWT		✓
29	6	56636-1		✓
30	7	56638-15		✓
31	8	~~PE~~P4045 Pest 50 yy 5.31.06		✓ Heptachlor ↓ 64% #1
32	9	P3982 C/T 0.5/1.0		✓
33	10	P4021 16/60 0.5		✓
34	99	P3862 PEM		✓
35	11	B053O6PAS, Florasil		✓ yy 6.1.06
36	12	L053O6PAST,		✓ endrin ketone marg ↑ #2 dw#1
37	13	L053O6PAST, ↓		✓ ↓
38	14	56625-1, 1:10, Florasil	did not report, used straight run	✓
39	15	-1, MS, 1:10		~~some failures~~ yy 6.1.06
40	16	-1, MSD, 1:10 ↓		✓ ↓
41	8	P4045 Pest 50		✓
42	9	P3982 C/T 0.5/1.0		✓
43	10	P4021 16/60 0.5		✓
44	99	P3862 PEM		✓
45	17	56625-1, Florasil		✓
46	18	-1, MS ↓		✓ some failures

Standard Dilutions: 1:2- 500 uL sample +500 uL hexane
1:5- 200 uL sample +800 uL hexane
Hexane Lot : 460160 1:10- 100 uL sample + 900 uL hexane
1:20- 50 uL sample + 950 uL hexane
MTBE Lot: _____ 1:100- 10 uLsample +990 uL hexane
1:1000- 1 uLsample +1000 uL hexane

C:\HPCHEM\2\METHODS\PBQ05166.M

Level 1 File Name	F55861.D
Level 2 File Name	F55862.D
Level 3 File Name	F55863.D
Level 4 File Name	F55864.D
Level 5 File Name	F55865.D
Level 6 File Name	F55866.D
Level 7 File Name	F55867.D

5-17-06 5.17.06

RESPONSE = X + YA + Z(A*A)

Name	R^2	R	X	Y	Z
TCX	1.000	1.000	1.04E+04	3.47E+07	0.00E-01
Ar1016 A	0.999	0.999	2.45E+04	1.65E+06	0.00E-01
Ar1016 B	0.999	1.000	3.80E+04	3.18E+06	0.00E-01
Ar1016 C	1.000	1.000	1.15E+04	1.34E+06	0.00E-01
Ar1260 A	0.999	1.000	2.75E+04	2.04E+06	0.00E-01
Ar1260 B	1.000	1.000	2.73E+04	2.45E+06	0.00E-01
Ar1260 C	1.000	1.000	8.21E+03	1.96E+06	0.00E-01
Ar1260 D	1.000	1.000	7.55E+03	1.93E+06	0.00E-01
DCB	0.999	0.999	9.77E+04	2.91E+07	0.00E-01
TCX #2	0.999	1.000	6.78E+04	3.70E+07	0.00E-01
Ar1016 A #2	0.997	0.999	4.34E+04	1.69E+06	0.00E-01
Ar1016 B #2	0.998	0.999	8.89E+04	3.83E+06	0.00E-01
Ar1016 C #2	0.997	0.999	3.37E+04	1.45E+06	0.00E-01
Ar1260 A #2	0.995	0.998	8.05E+04	2.28E+06	0.00E-01
Ar1260 B #2	0.996	0.998	1.10E+05	3.42E+06	0.00E-01
Ar1260 C #2	0.997	0.999	6.09E+04	2.54E+06	0.00E-01
Ar1260 D #2	0.999	0.999	7.76E+04	5.97E+06	0.00E-01
DCB #2	0.994	0.997	2.64E+05	3.87E+07	0.00E-01

Quantitation Report (QT Reviewed)

Signal #1 : C:\HPCHEM\2\DATA\051606-F\F55861.D\ECD1A.CH Vial: 13
Signal #2 : C:\HPCHEM\2\DATA\051606-F\F55861.D\ECD2B.CH
Acq On : 16 May 2006 7:38 pm Operator:
Sample : P4016 16/60 0.01 Inst : Instrumen
Misc : Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 17 9:56 2006 Quant Results File: PBQ05166.RES

Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 09:54:21 2006
Response via : Initial Calibration
DataAcq Meth : PSTQUICK.M

5-17-06 5-17-06

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film

Compound	RT#1	RT#2	Resp#1	Resp#2	ug/ml	ug/ml
System Monitoring Compounds						
1) S TCX	2.64	2.72	61307	77106	0.002	0.001 #
Spiked Amount 0.100	Range 46 - 122		Recovery =		2.00%#	1.00%#
9) S DCB	10.39	10.04	79033	133576	N.D. m	N.D. m
Spiked Amount 0.100	Range 26 - 86		Recovery =		0.00%#	0.00%#
Target Compounds						
2) L1 Ar1016 A	3.43	3.37	19531	27177	N.D. m	N.D.
3) L1 Ar1016 B	3.92	3.86	43320	57040	0.001	N.D. #
4) L1 Ar1016 C	4.09	4.03	15699	21986	0.004	N.D. #
Sum Ar1016 A			59019	0	0.001	N.D.
Average Ar1016 A					0.000	0.000
5) L2 Ar1260 A	6.67	6.35	30275	39717	0.002	N.D. m#
6) L2 Ar1260 B	7.03	6.81	33089	60122	0.001	N.D. m#
7) L2 Ar1260 C	7.92	7.53	22273	35850	0.006	N.D. m#
8) L2 Ar1260 D	8.49	8.59	22028	68293	0.004	N.D. #
Sum Ar1260 A			107665	0	0.013	N.D.
Average Ar1260 A					0.003	0.000

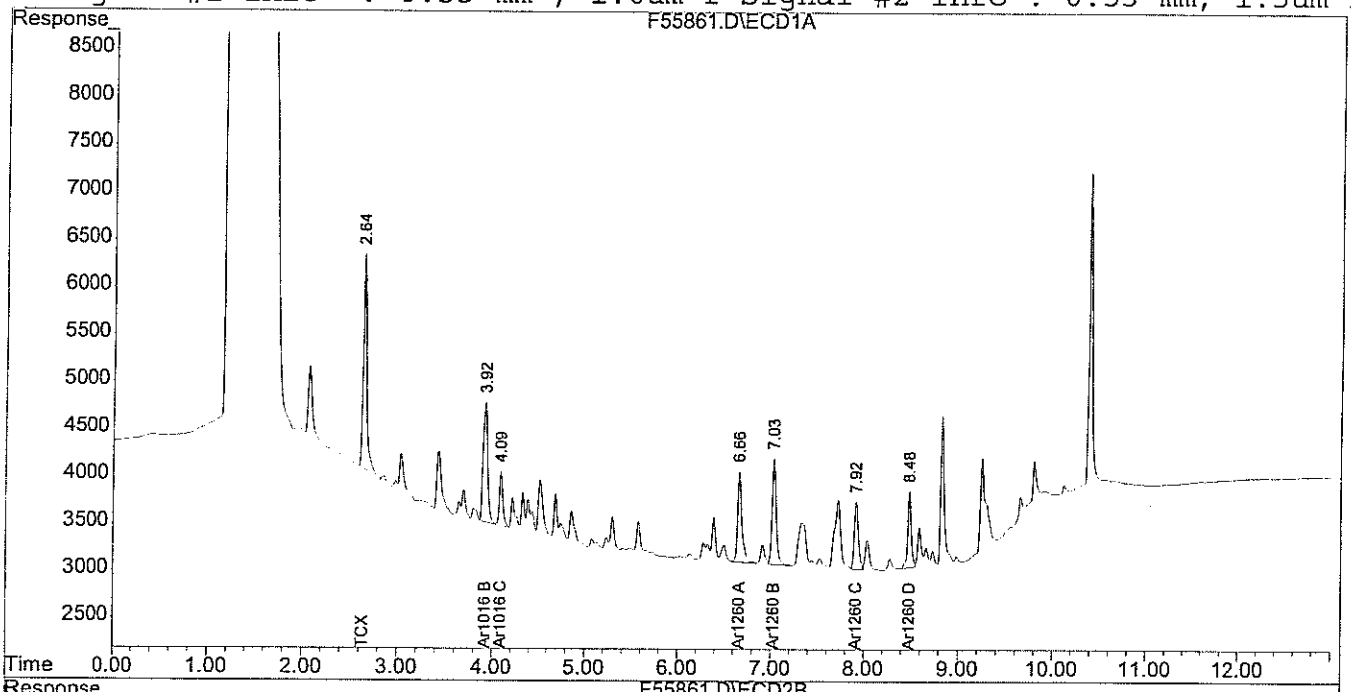
(f)=RT Delta > 1/2 Window (#)=Amounts differ by > 25% (m)=manual int.

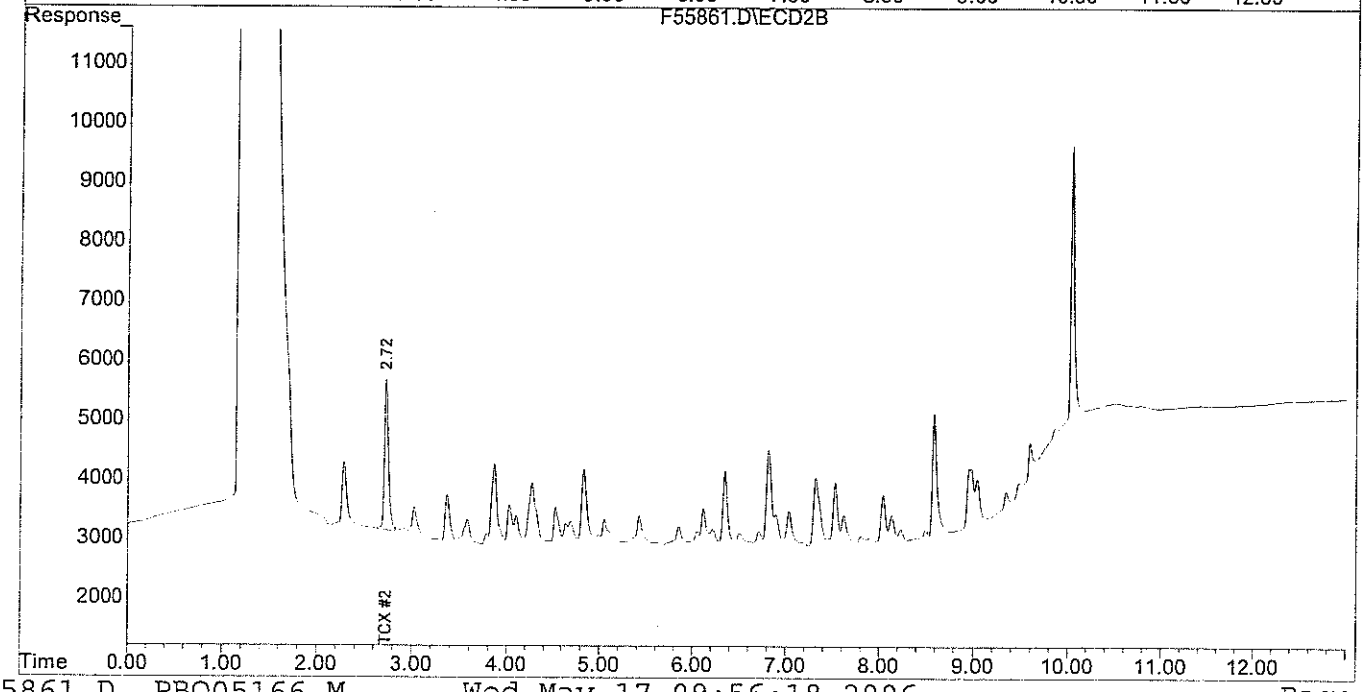
Quantitation Report (QT Reviewed)

Signal #1 : C:\HPCHEM\2\DATA\051606-F\F55861.D\ECD1A.CH Vial: 13
Signal #2 : C:\HPCHEM\2\DATA\051606-F\F55861.D\ECD2B.CH
Acq On : 16 May 2006 7:38 pm Operator:
Sample : P4016 16/60 0.01 Inst : Instrumen
Misc : Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 17 9:56 2006 Quant Results File: PBQ05166.RES

Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 09:54:21 2006
Response via : Multiple Level Calibration
DataAcq Meth : PSTQUICK.M

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film



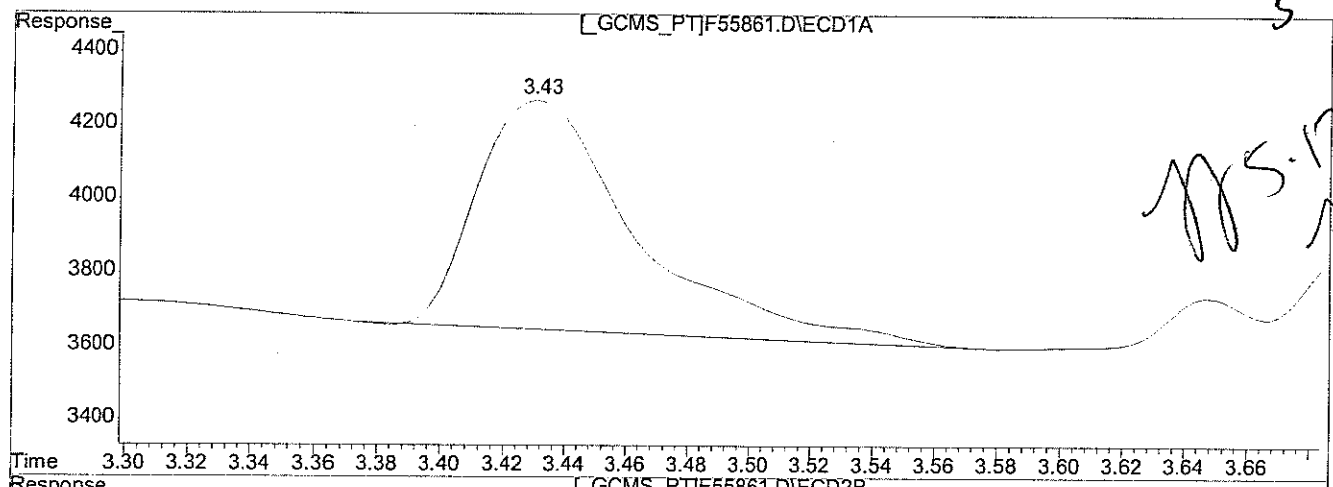


Quantitation Report (Qedit)

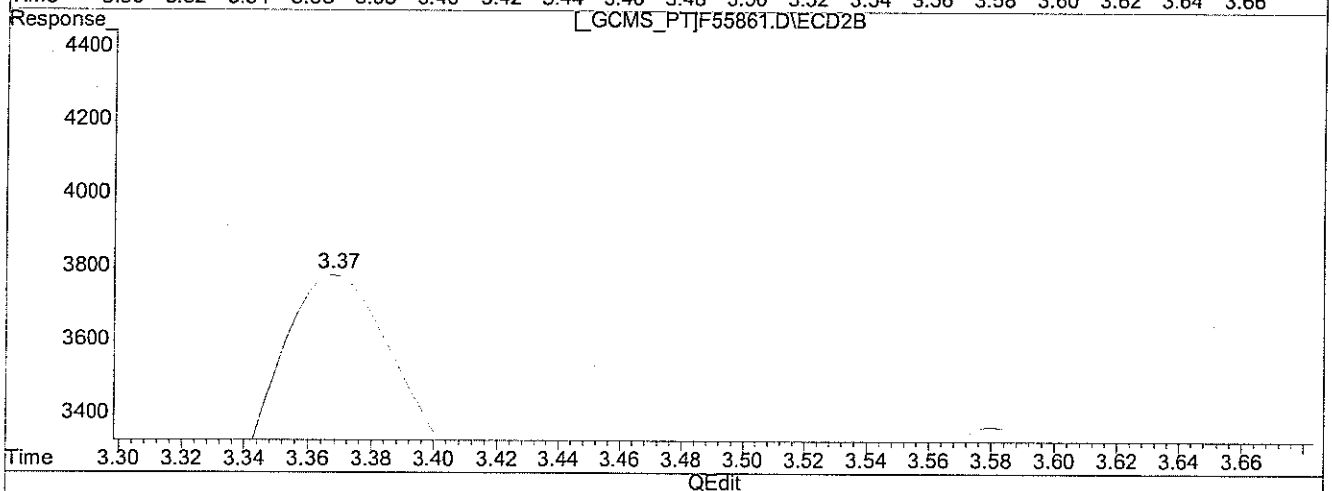
Signal #1 : C:\HPCHEM\2\DATA\051606-F\F55861.D\ECD1A.CH Vial: 13
Signal #2 : C:\HPCHEM\2\DATA\051606-F\F55861.D\ECD2B.CH
Acq On : 16 May 2006 7:38 pm Operator:
Sample : P4016 16/60 0.01 Inst : Instrumen
Misc : Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 17 9:54 2006 Quant Results File: PBQ05166.RES

Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 09:54:21 2006
Response via : Multiple Level Calibration

5.17.06



5.17.06 Shoulder



(2) Ar1016 A (L1)

R.T.	Response	Conc
3.43	22374	-0.00
3.92	43320	0.00
4.09	15699	0.00

(2) Ar1016 A #2 (L1)

R.T.	Response	Conc
3.37	27177	-0.01
3.86	57040	-0.01
4.03	21986	-0.01

(+) = Expected Retention Time

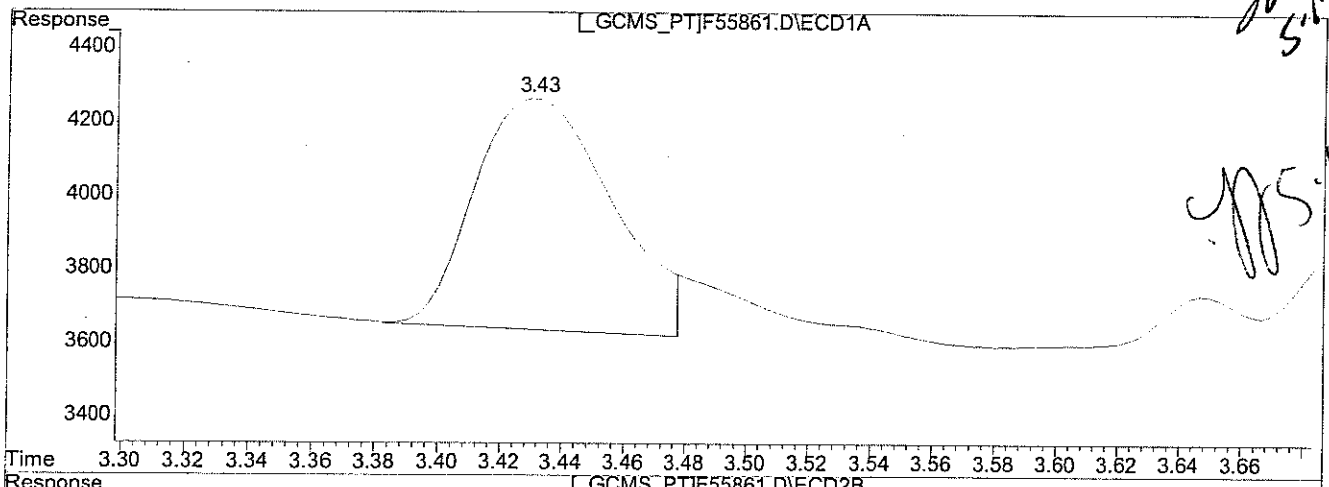
Quantitation Report (Qedit)

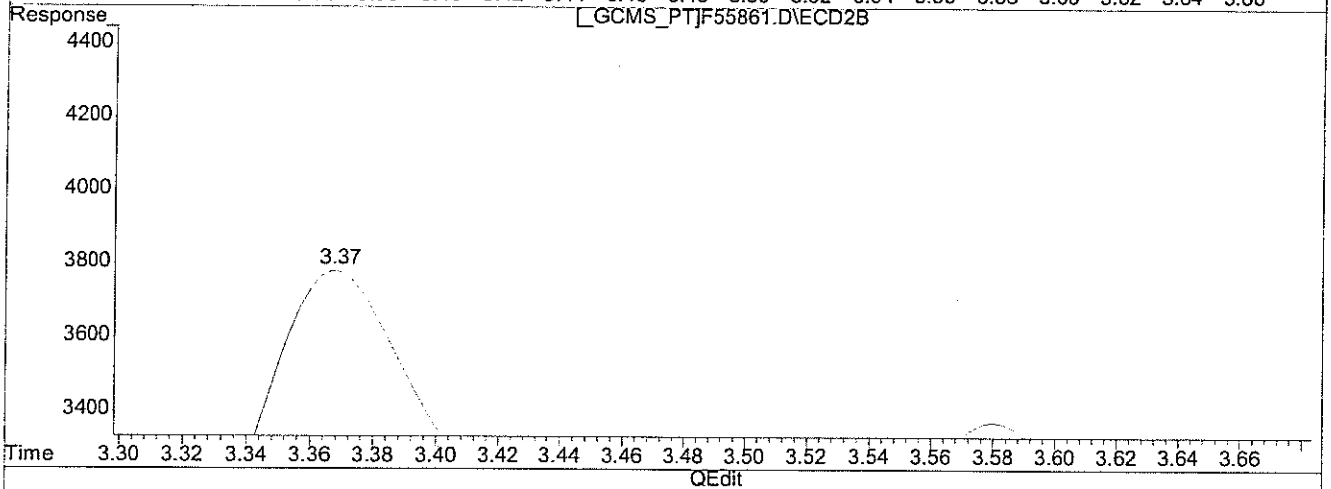
Signal #1 : C:\HPCHEM\2\DATA\051606-F\F55861.D\ECD1A.CH Vial: 13
Signal #2 : C:\HPCHEM\2\DATA\051606-F\F55861.D\ECD2B.CH
Acq On : 16 May 2006 7:38 pm Operator:
Sample : P4016 16/60 0.01 Inst : Instrumen
Misc : Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 17 9:54 2006 Quant Results File: PBQ05166.RES

Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 09:54:21 2006
Response via : Multiple Level Calibration

5.17.06

5.17.06





(2) Ar1016 A #2 (L1)

R.T.	Response	Conc
3.43	19531	-0.00
3.92	43320	0.00
4.09	15699	0.00

(2) Ar1016 A #2 (L1)

R.T.	Response	Conc
3.37	27177	-0.01
3.86	57040	-0.01
4.03	21986	-0.01

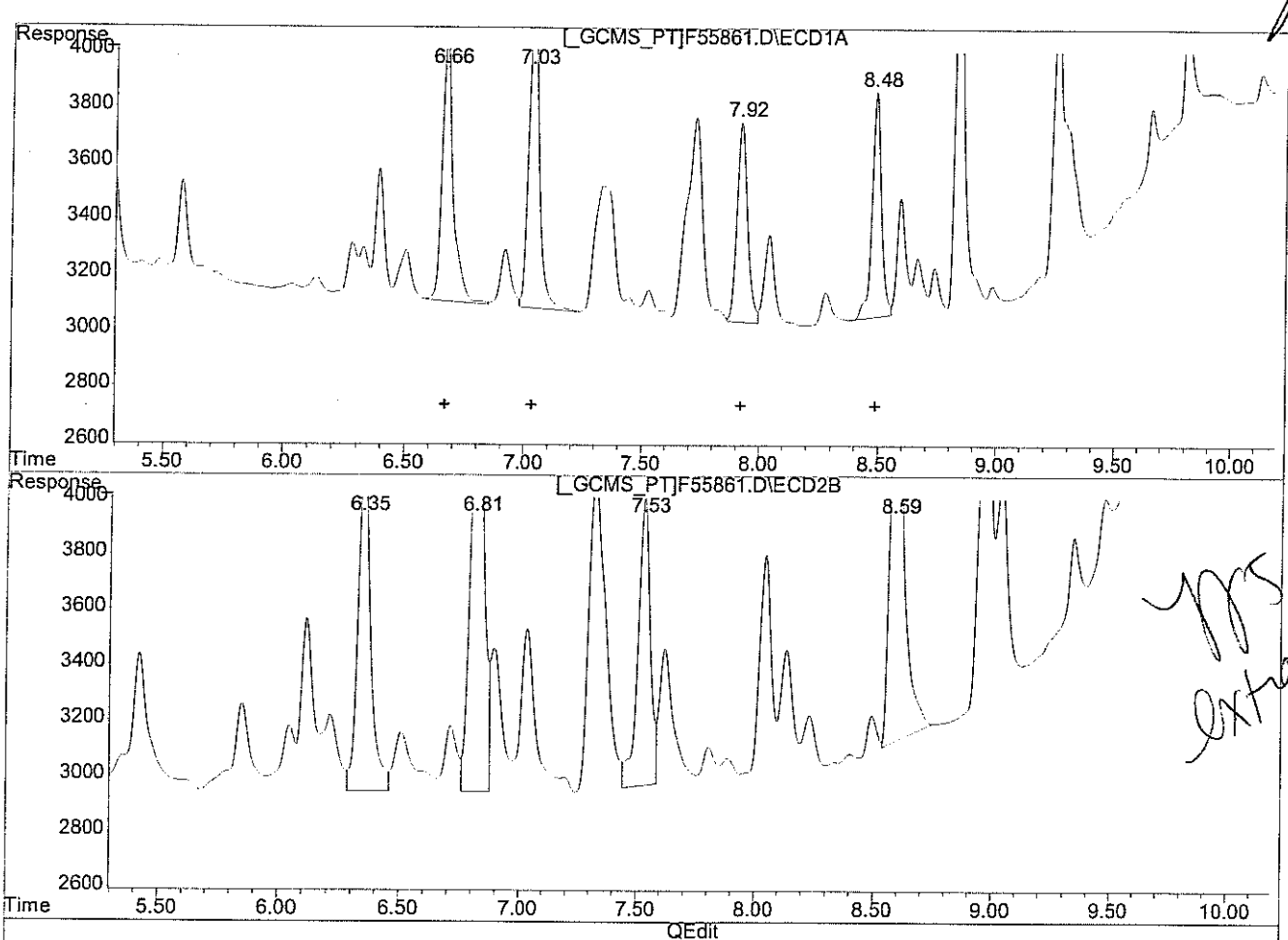
(+) = Expected Retention Time

Quantitation Report (Qedit)

Signal #1 : C:\HPCHEM\2\DATA\051606-F\F55861.D\ECD1A.CH Vial: 13
Signal #2 : C:\HPCHEM\2\DATA\051606-F\F55861.D\ECD2B.CH
Acq On : 16 May 2006 7:38 pm Operator:
Sample : P4016 16/60 0.01 Inst : Instrumen
Misc : Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 17 9:54 2006 Quant Results File: PBQ05166.RES

Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 09:54:21 2006
Response via : Multiple Level Calibration

5.17.06



5.17.06
Extraction

(5) Ar1260 A (L2)

R.T.	Response	Conc
6.67	30275	0.00
7.03	33089	0.00
7.92	22273	0.01
8.49	22028	0.00

(5) Ar1260 A #2 (L2)

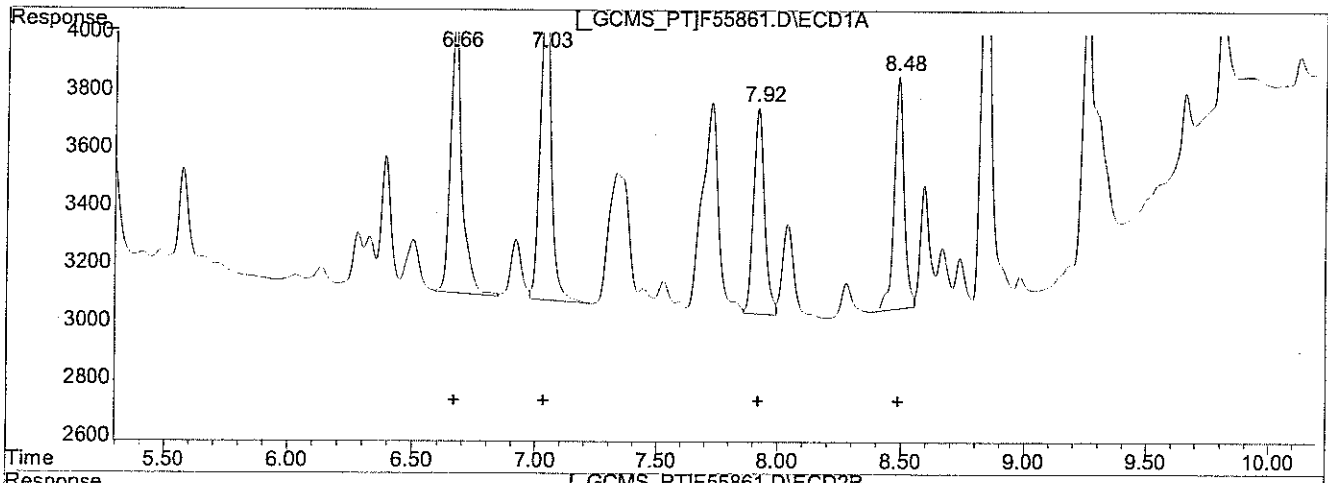
R.T.	Response	Conc
6.35	45036	-0.02
6.82	61803	-0.01
7.53	37673	-0.01
8.59	68293	-0.00

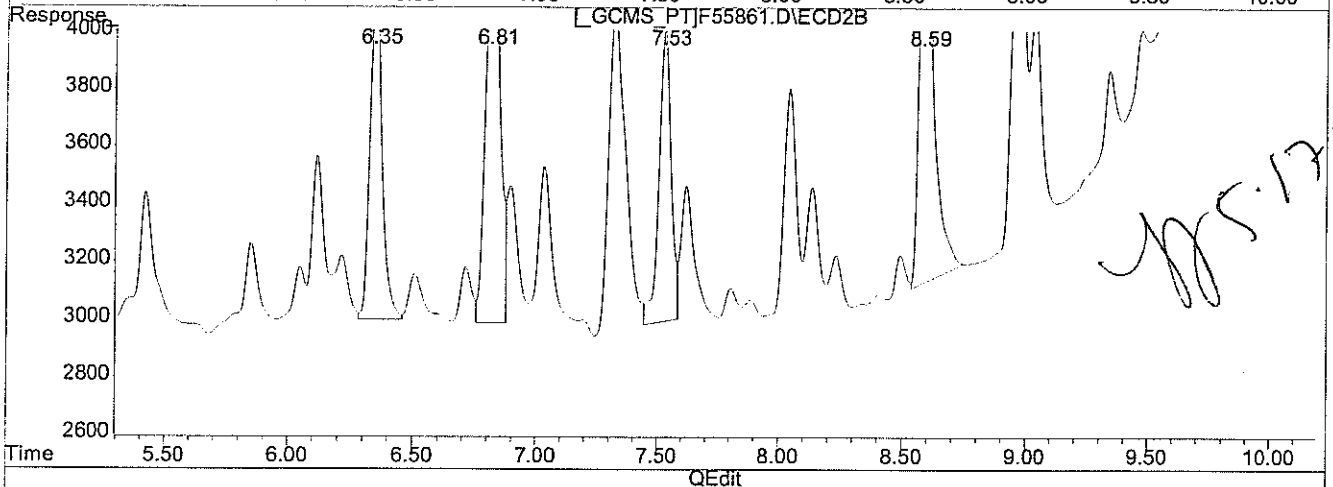
(+) = Expected Retention Time

Quantitation Report (Qedit)

Signal #1 : C:\HPCHEM\2\DATA\051606-F\F55861.D\ECD1A.CH Vial: 13
Signal #2 : C:\HPCHEM\2\DATA\051606-F\F55861.D\ECD2B.CH
Acq On : 16 May 2006 7:38 pm Operator:
Sample : P4016 16/60 0.01 Inst : Instrumen
Misc : Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 17 9:54 2006 Quant Results File: PBQ05166.RES

Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 09:54:21 2006
Response via : Multiple Level Calibration





QEdit

(5) Ar1260 A #2 (L2)

R.T.	Response	Conc
6.67	30275	0.00
7.03	33089	0.00
7.92	22273	0.01
8.49	22028	0.00

(5) Ar1260 A #2 (L2)

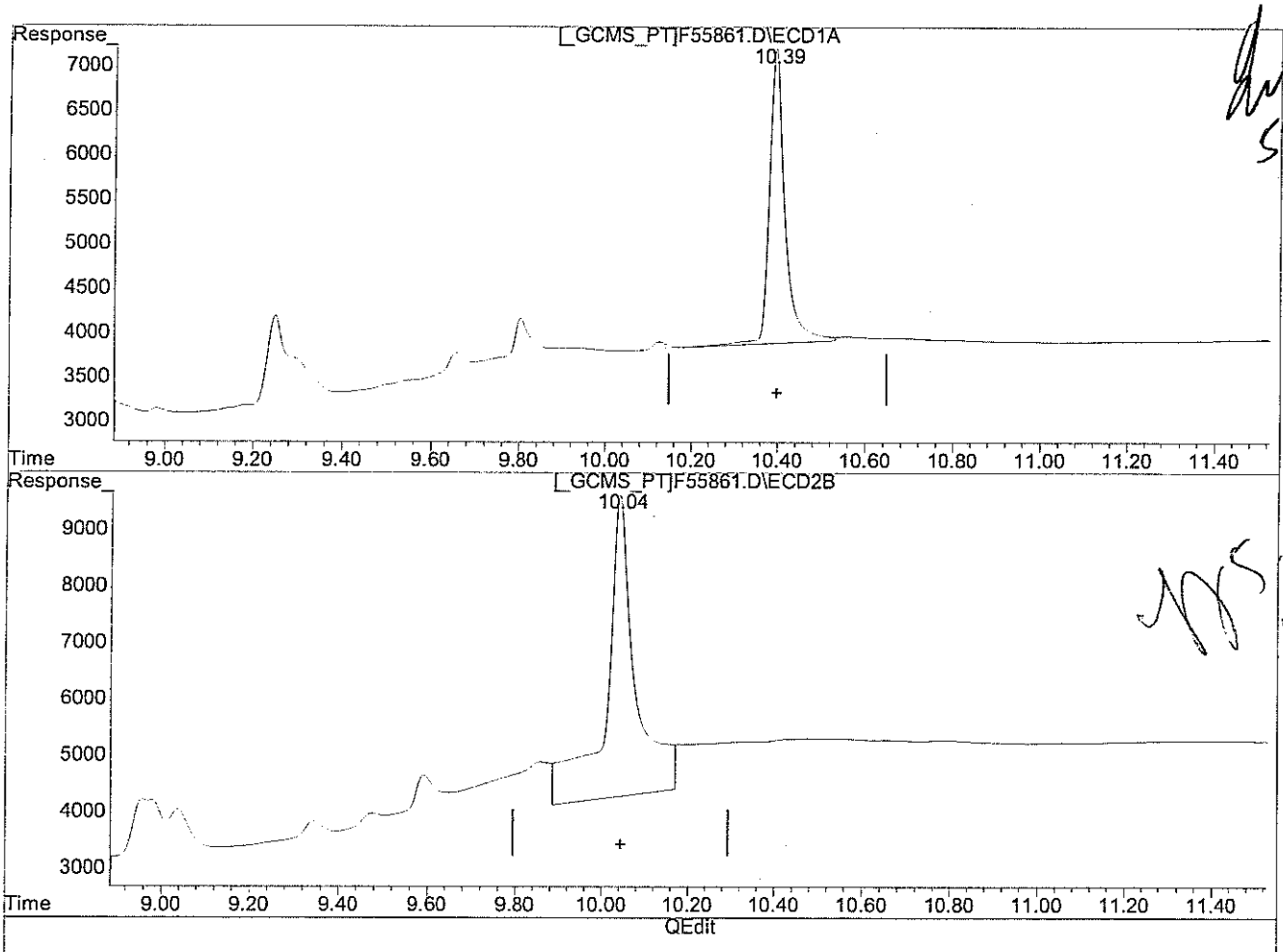
R.T.	Response	Conc
6.35	39717	-0.02
6.81	60122	-0.01
7.53	35850	-0.01
8.59	68293	-0.00

(+) = Expected Retention Time

Quantitation Report (Qedit)

Signal #1 : C:\HPCHEM\2\DATA\051606-F\F55861.D\ECD1A.CH Vial: 13
Signal #2 : C:\HPCHEM\2\DATA\051606-F\F55861.D\ECD2B.CH
Acq On : 16 May 2006 7:38 pm Operator:
Sample : P4016 16/60 0.01 Inst : Instrumen
Misc : Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 17 9:54 2006 Quant Results File: PBQ05166.RES

Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 09:54:21 2006
Response via : Multiple Level Calibration



(9) DCB (S)

10.39min -0.001ug/ml

response 81717

(9) DCB #2 (S)

10.04min -0.000ug/ml

response 257756

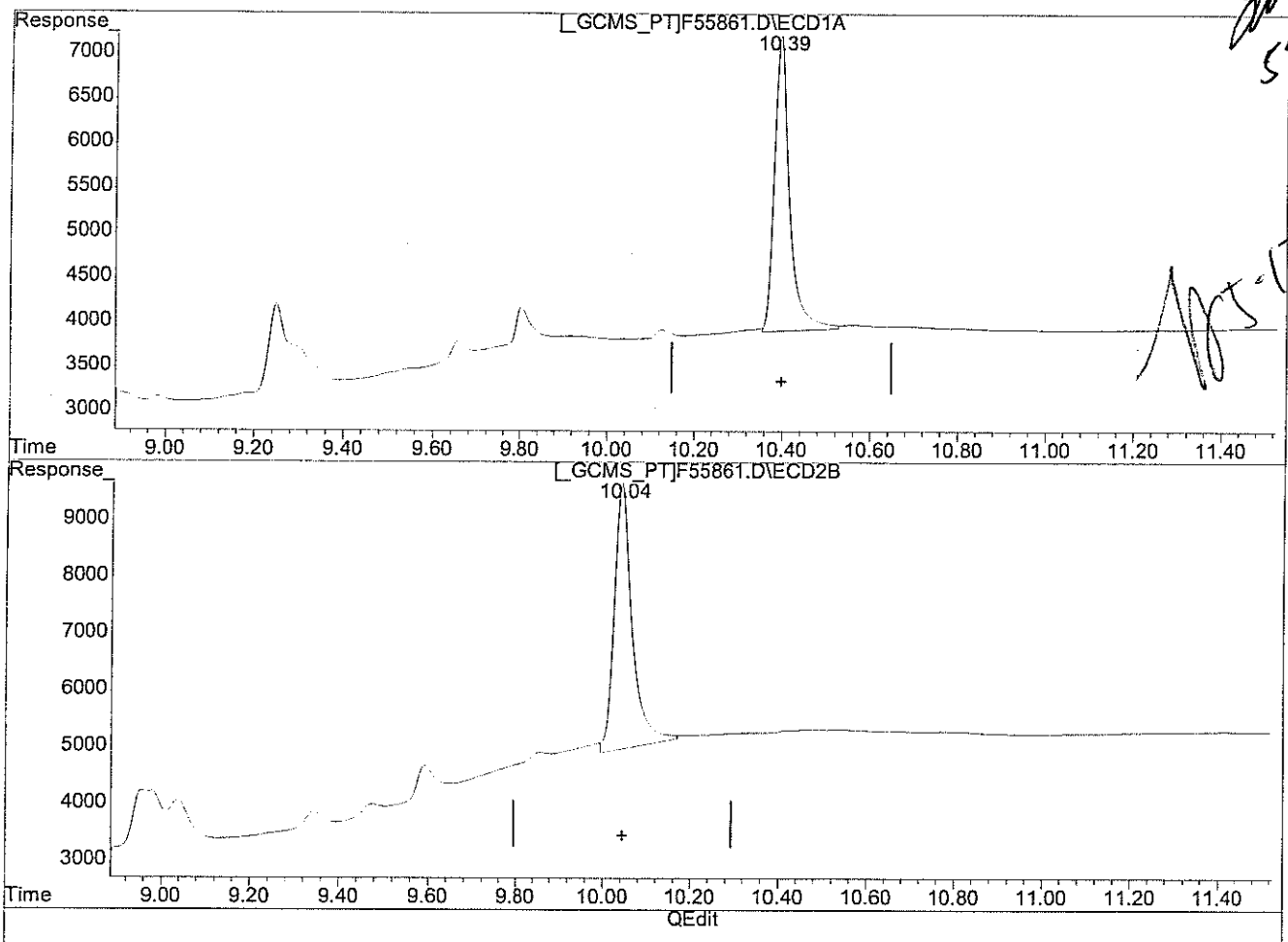
(+) = Expected Retention Time

F55861.D PBQ05166.M Wed May 17 09:56:02 2006

Quantitation Report (Qedit)

Signal #1 : C:\HPCHEM\2\DATA\051606-F\F55861.D\ECD1A.CH Vial: 13
Signal #2 : C:\HPCHEM\2\DATA\051606-F\F55861.D\ECD2B.CH
Acq On : 16 May 2006 7:38 pm Operator:
Sample : P4016 16/60 0.01 Inst : Instrumen
Misc : Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 17 9:54 2006 Quant Results File: PBQ05166.RES

Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 09:54:21 2006
Response via : Multiple Level Calibration



(9) DCB (S)

10.39min -0.001ug/ml m

response 79033

(9) DCB #2 (S)

10.04min -0.003ug/ml m

response 133576

(+) = Expected Retention Time

F55861.D PBQ05166.M Wed May 17 09:56:13 2006

Quantitation Report (QT Reviewed)

Signal #1 : C:\HPCHEM\2\DATA\051606-F\F55862.D\ECD1A.CH Vial: 14
Signal #2 : C:\HPCHEM\2\DATA\051606-F\F55862.D\ECD2B.CH
Acq On : 16 May 2006 7:53 pm Operator:
Sample : P4017 16/60 0.02 Inst : Instrumen
Misc : Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 17 9:56 2006 Quant Results File: PBQ05166.RES

Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 09:54:21 2006
Response via : Initial Calibration
DataAcq Meth : PSTQUICK.M

5-17-06 5-17-06

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film

Compound	RT#1	RT#2	Resp#1	Resp#2	ug/ml	ug/ml
System Monitoring Compounds						
1) S TCX	2.64	2.73	122756	153548	0.004	0.003
Spiked Amount 0.100	Range 46 - 122			Recovery =	4.00%#	3.00%#
9) S DCB	10.40	10.04	155614	248581	0.001m	N.D. m#
Spiked Amount 0.100	Range 26 - 86			Recovery =	1.00%#	0.00%#
Target Compounds						
2) L1 Ar1016 A	3.44	3.37	42784	51825	0.011	0.006 #
3) L1 Ar1016 B	3.93	3.87	80946	111881	0.014	0.008 #
4) L1 Ar1016 C	4.09	4.03	29570	41472	0.016	0.008 #
Sum Ar1016 A			153300	205178	0.041	0.022
Average Ar1016 A					0.014	0.007
5) L2 Ar1260 A	6.67	6.35	53627	81296	0.014	0.002 #
6) L2 Ar1260 B	7.04	6.82	59571	113223	0.013	0.004 #
7) L2 Ar1260 C	7.92	7.54	42912	71915	0.017	0.007 #
8) L2 Ar1260 D	8.49	8.59	42372	139041	0.016	0.012 #
Sum Ar1260 A			198483	405475	0.060	0.024
Average Ar1260 A					0.015	0.006

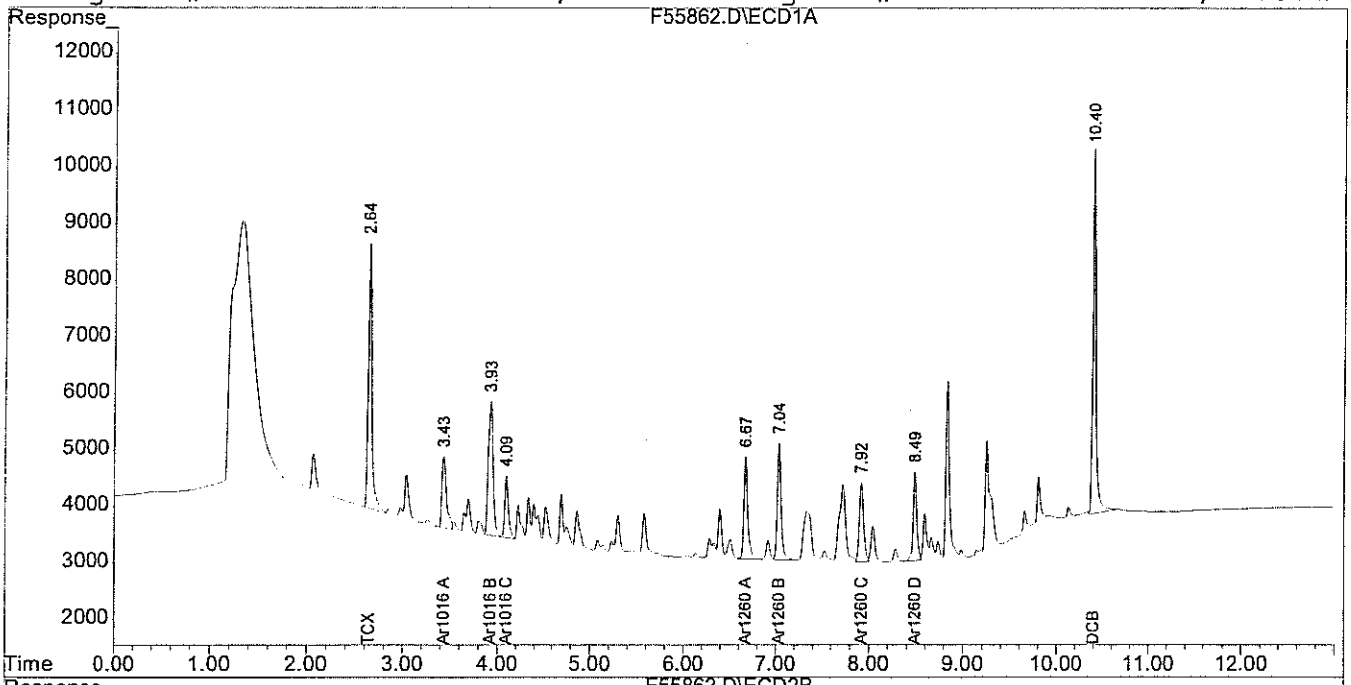
(f)=RT Delta > 1/2 Window (#)=Amounts differ by > 25% (m)=manual int.

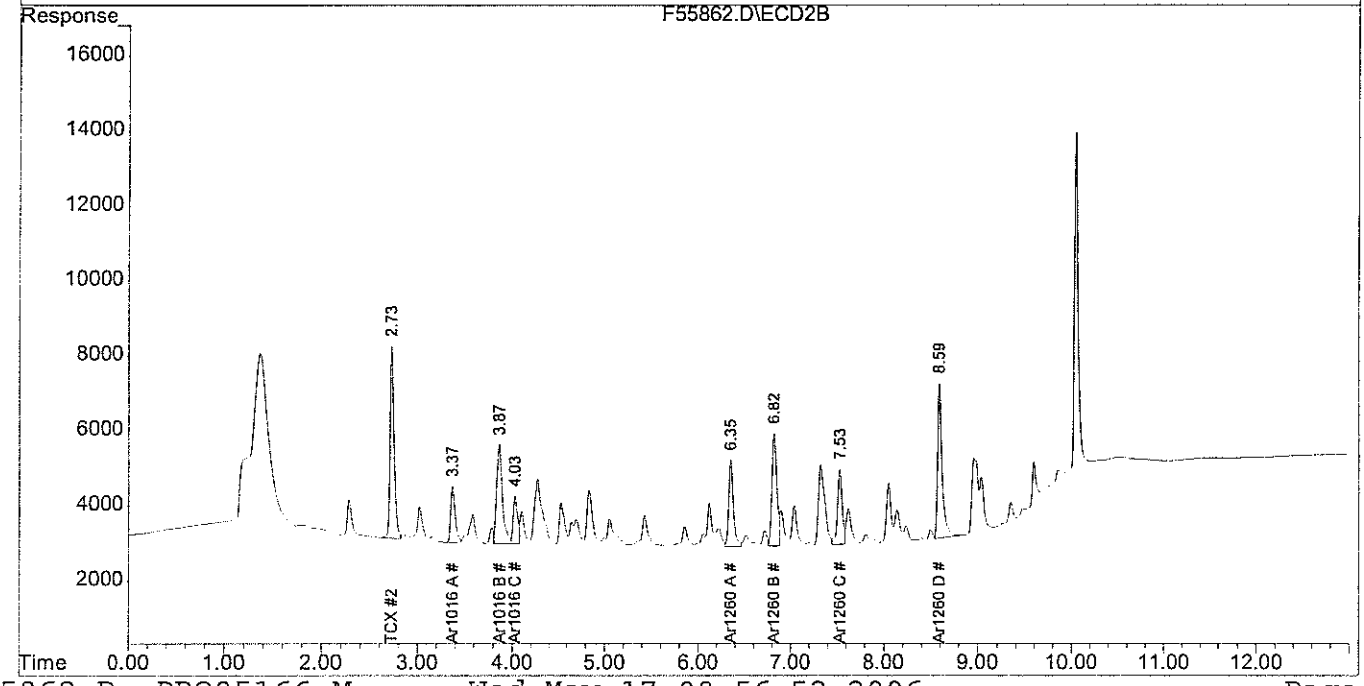
Quantitation Report (QT Reviewed)

Signal #1 : C:\HPCHEM\2\DATA\051606-F\F55862.D\ECD1A.CH Vial: 14
Signal #2 : C:\HPCHEM\2\DATA\051606-F\F55862.D\ECD2B.CH
Acq On : 16 May 2006 7:53 pm Operator:
Sample : P4017 16/60 0.02 Inst : Instrumen
Misc : Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 17 9:56 2006 Quant Results File: PBQ05166.RES

Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 09:54:21 2006
Response via : Multiple Level Calibration
DataAcq Meth : PSTQUICK.M

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film

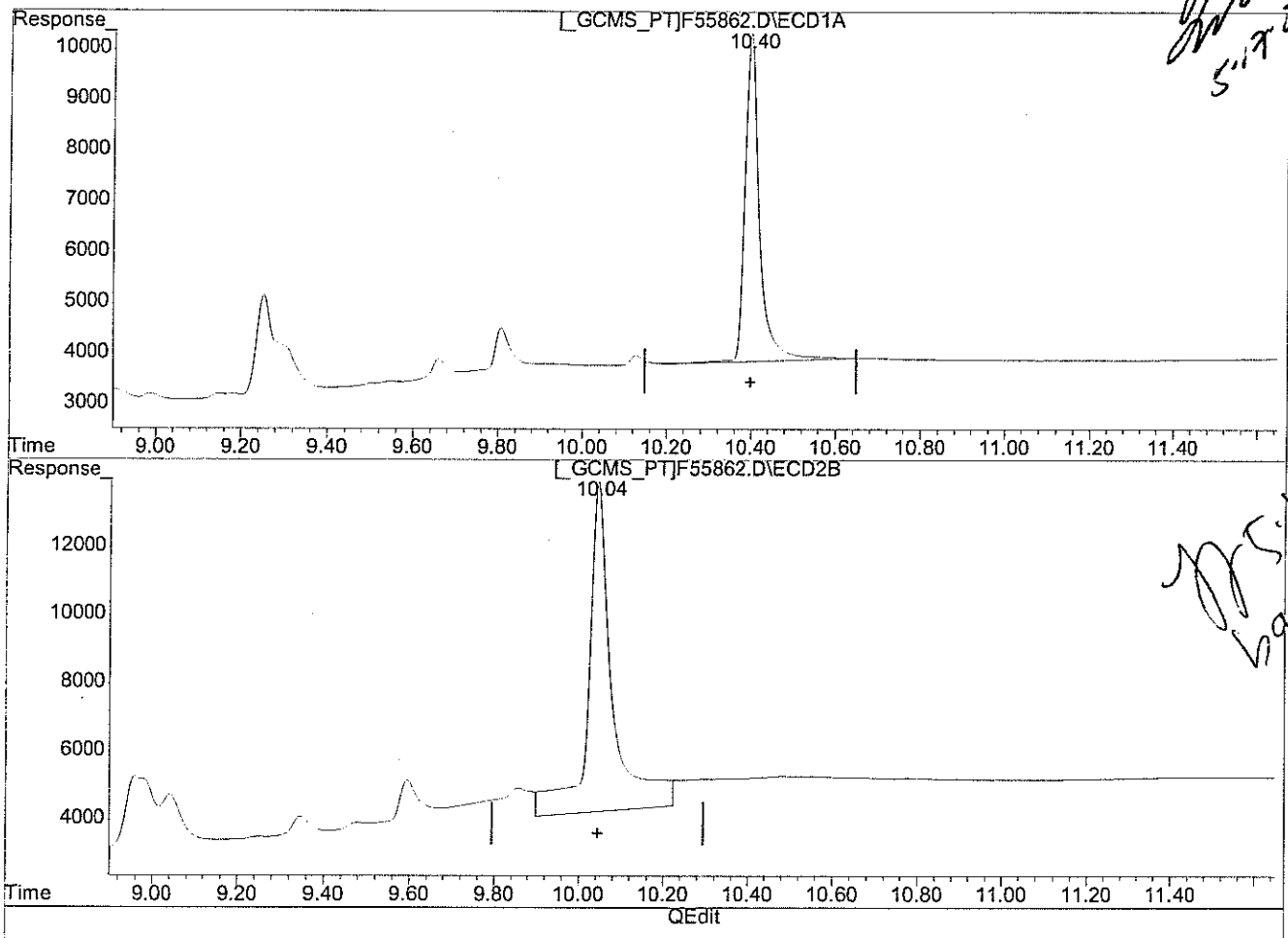




Quantitation Report (Qedit)

Signal #1 : C:\HPCHEM\2\DATA\051606-F\F55862.D\ECD1A.CH Vial: 14
Signal #2 : C:\HPCHEM\2\DATA\051606-F\F55862.D\ECD2B.CH
Acq On : 16 May 2006 7:53 pm Operator:
Sample : P4017 16/60 0.02 Inst : Instrumen
Misc : Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 17 9:56 2006 Quant Results File: PBQ05166.RES

Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 09:54:21 2006
Response via : Multiple Level Calibration



QEdit

(9) DCB (S)

10.40min 0.002ug/ml

response 156938

(9) DCB #2 (S)

10.04min 0.003ug/ml

response 389832

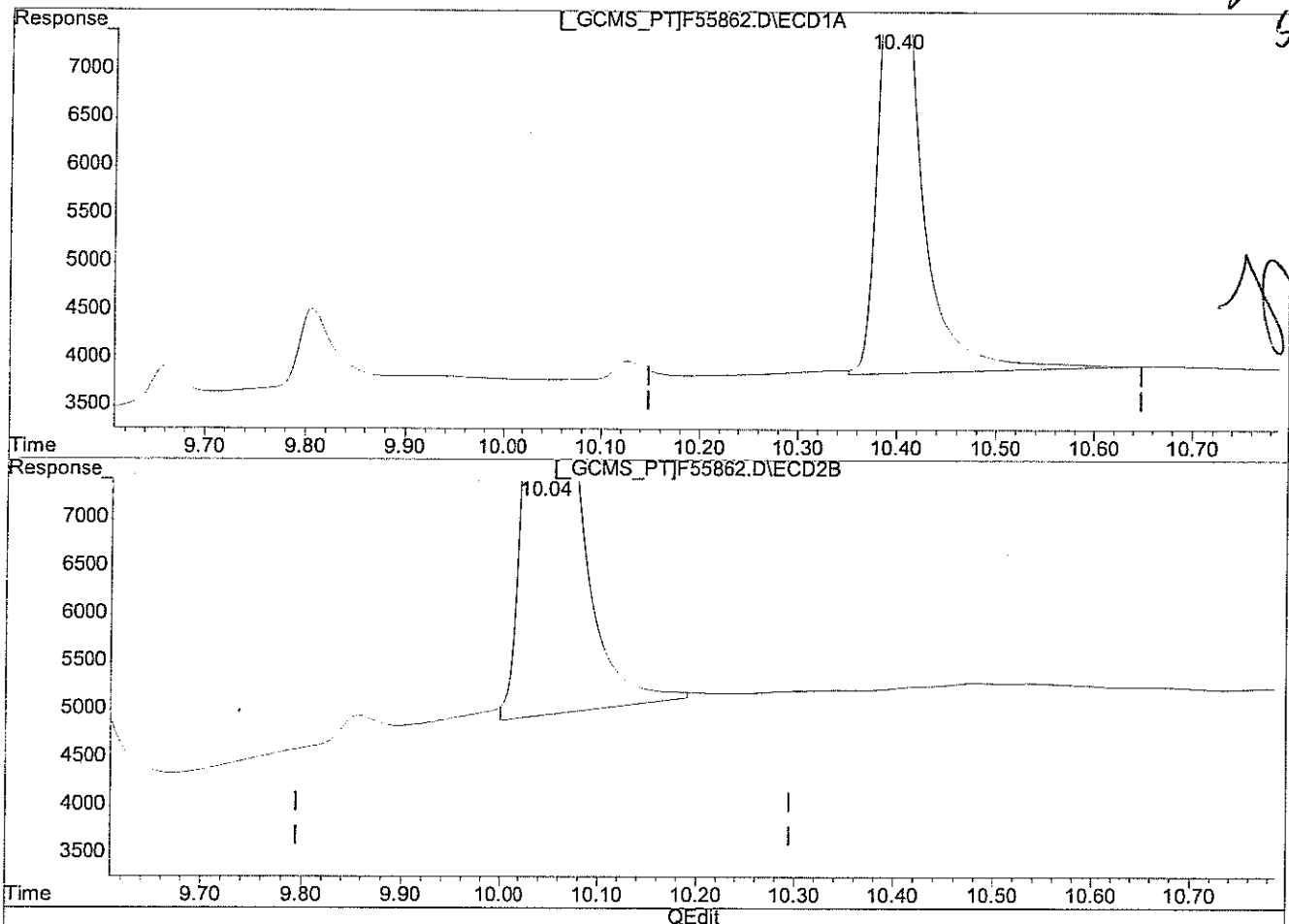
(+) = Expected Retention Time

Quantitation Report (Qedit)

Signal #1 : C:\HPCHEM\2\DATA\051606-F\F55862.D\ECD1A.CH Vial: 14
Signal #2 : C:\HPCHEM\2\DATA\051606-F\F55862.D\ECD2B.CH
Acq On : 16 May 2006 7:53 pm Operator:
Sample : P4017 16/60 0.02 Inst : Instrumen
Misc : Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 17 9:56 2006 Quant Results File: PBQ05166.RES

Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 09:54:21 2006
Response via : Multiple Level Calibration

5.17.06



5.17.06

(9) DCB (S)

10.40min 0.001ug/ml m

response 155614

(9) DCB #2 (S)

10.04min -0.000ug/ml m

response 248581

(+) = Expected Retention Time

F55862.D PBQ05166.M Wed May 17 09:56:48 2006

Quantitation Report (QT Reviewed)

Signal #1 : C:\HPCHEM\2\DATA\051606-F\F55863.D\ECD1A.CH Vial: 15
Signal #2 : C:\HPCHEM\2\DATA\051606-F\F55863.D\ECD2B.CH
Acq On : 16 May 2006 8:07 pm Operator:
Sample : P4018 16/60 0.05 Inst : Instrumen
Misc : Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 17 10:04 2006 Quant Results File: PBQ05166.RES

Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 09:54:21 2006
Response via : Initial Calibration
DataAcq Meth : PSTQUICK.M

5.17.06 5.17.06

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film

Compound	RT#1	RT#2	Resp#1	Resp#2	ug/ml	ug/ml
System Monitoring Compounds						
1) S TCX	2.64	2.73	328762	401821	0.010	0.010
Spiked Amount 0.100	Range 46 - 122		Recovery =		10.00%#	10.00%#
9) S DCB	10.40	10.05	369688	589584	0.009m	0.009m
Spiked Amount 0.100	Range 26 - 86		Recovery =		9.00%#	9.00%#
Target Compounds						
2) L1 Ar1016 A	3.44	3.37	102888	121190	0.051	0.050
3) L1 Ar1016 B	3.93	3.87	187238	263561	0.051	0.050
4) L1 Ar1016 C	4.10	4.03	75426	99033	0.054	0.051
Sum Ar1016 A			365551	483784	0.156	0.151
Average Ar1016 A					0.052	0.050
5) L2 Ar1260 A	6.67	6.35	126226	190696	0.053	0.054
6) L2 Ar1260 B	7.04	6.82	148952	269972	0.053	0.054
7) L2 Ar1260 C	7.92	7.53	105396	183351	0.052	0.056
8) L2 Ar1260 D	8.49	8.59	102899	351344	0.051	0.051
Sum Ar1260 A			483473	995362	0.209	0.216
Average Ar1260 A					0.052	0.054

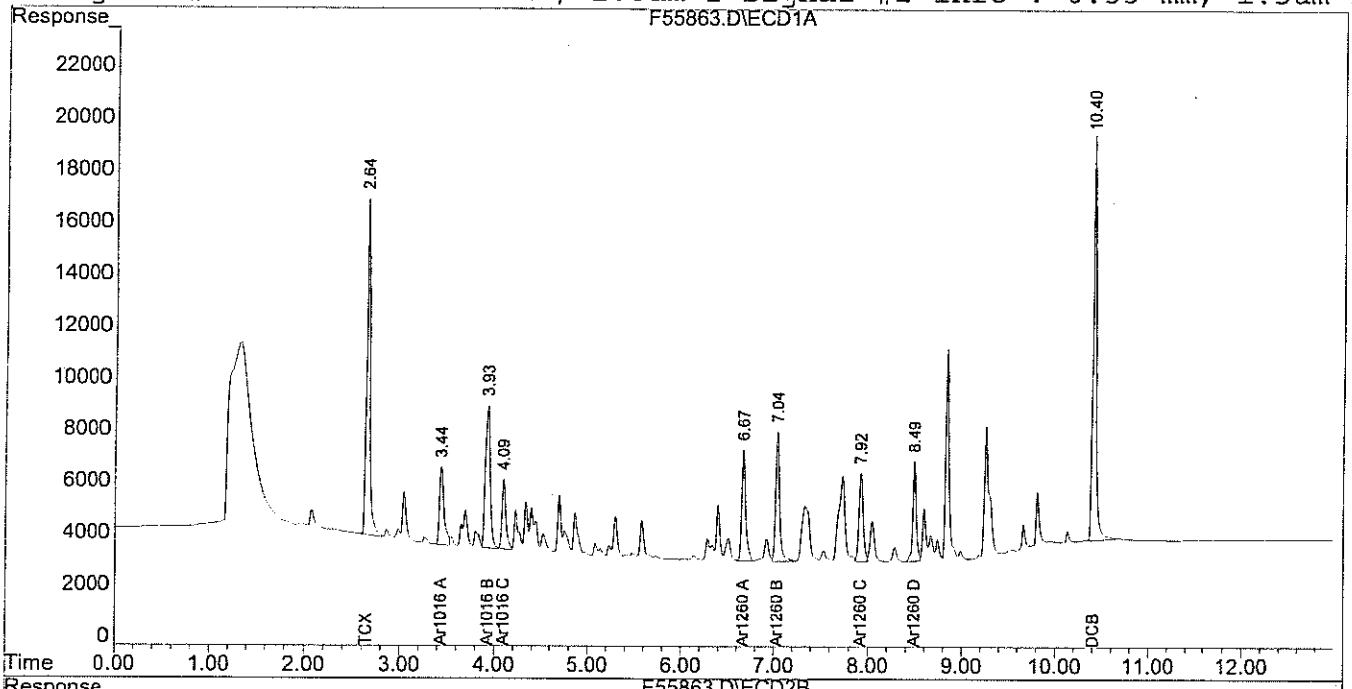
(f)=RT Delta > 1/2 Window (#)=Amounts differ by > 25% (m)=manual int.

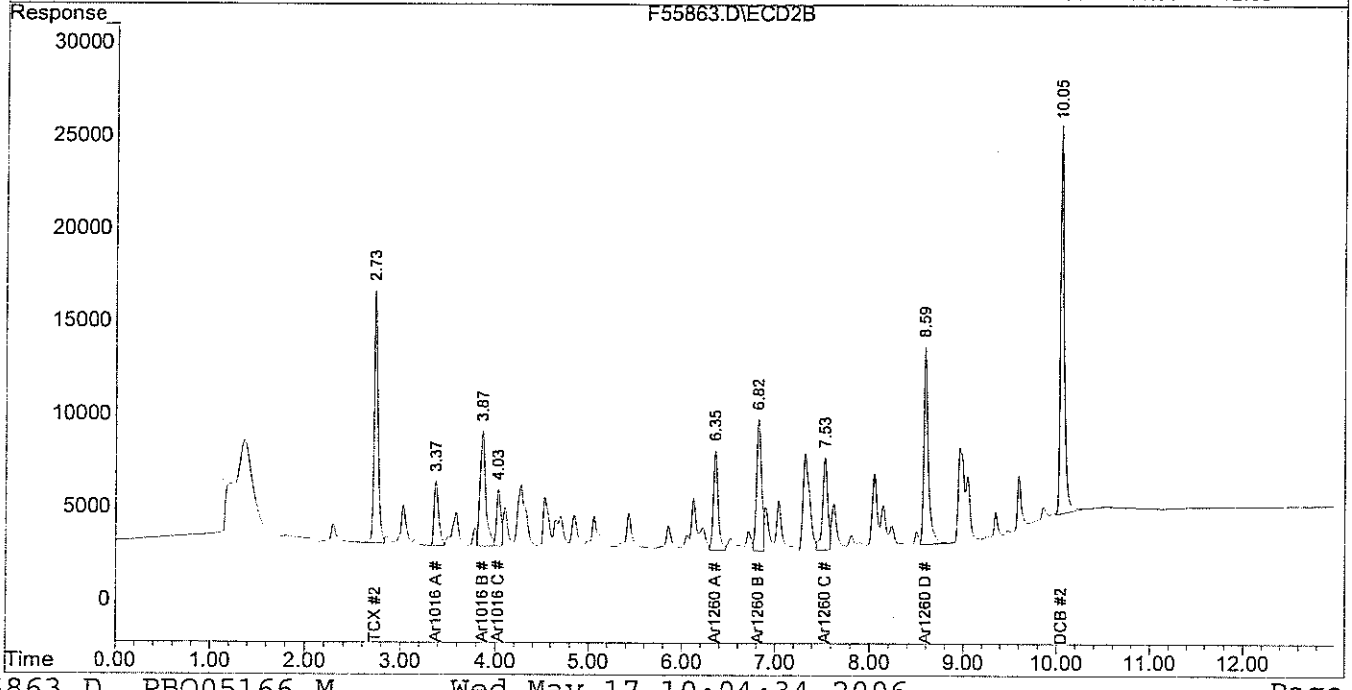
Quantitation Report (QT Reviewed)

Signal #1 : C:\HPCHEM\2\DATA\051606-F\F55863.D\ECD1A.CH Vial: 15
Signal #2 : C:\HPCHEM\2\DATA\051606-F\F55863.D\ECD2B.CH
Acq On : 16 May 2006 8:07 pm Operator:
Sample : P4018 16/60 0.05 Inst : Instrumen
Misc : Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 17 10:04 2006 Quant Results File: PBQ05166.RES

Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 09:54:21 2006
Response via : Multiple Level Calibration
DataAcq Meth : PSTQUICK.M

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film



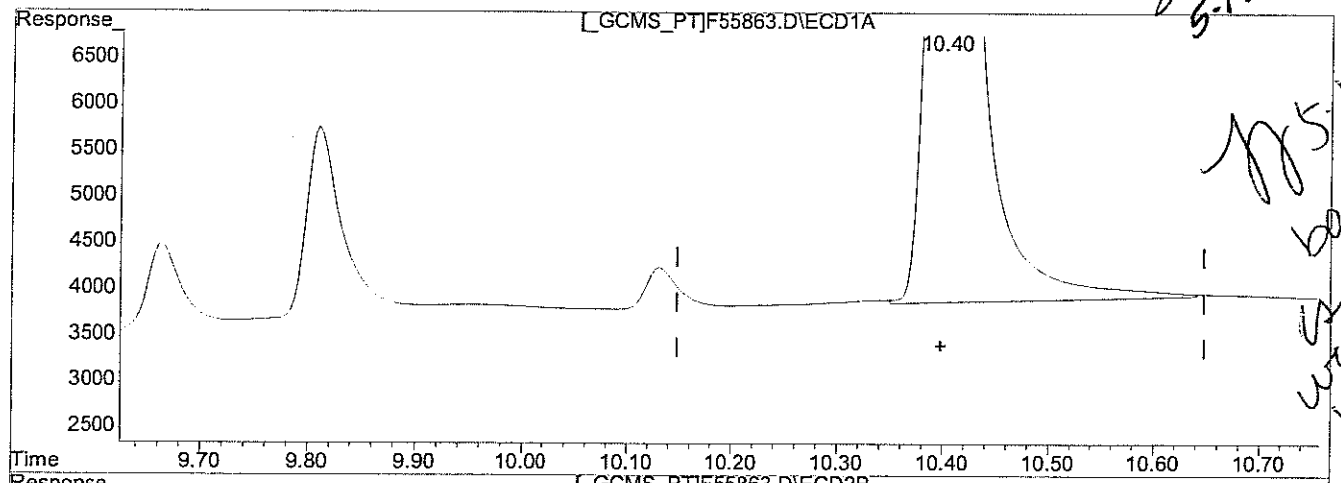


Quantitation Report (Qedit)

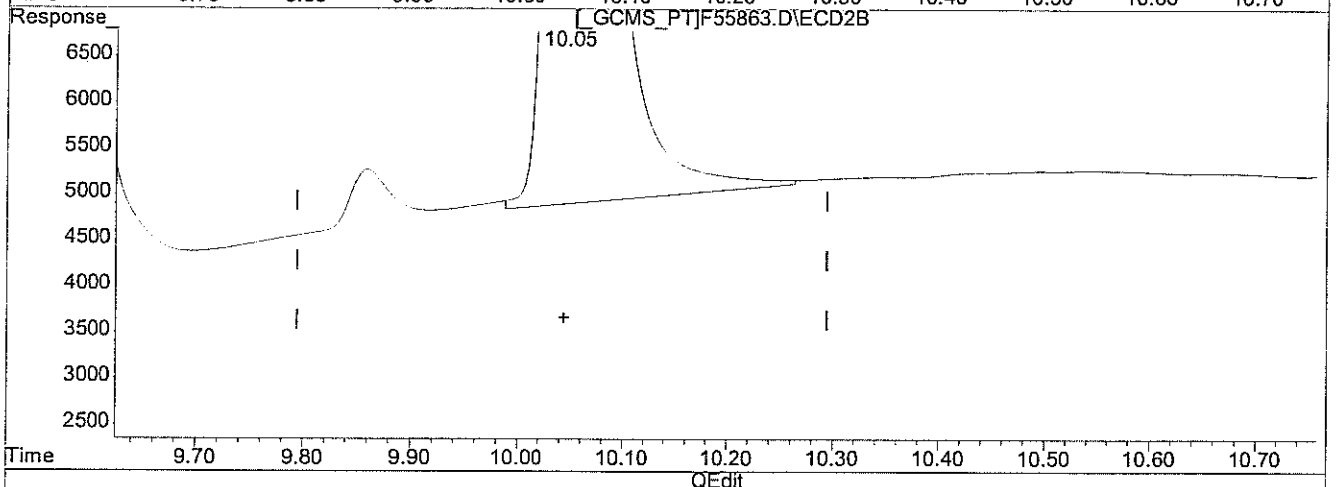
Signal #1 : C:\HPCHEM\2\DATA\051606-F\F55863.D\ECD1A.CH Vial: 15
Signal #2 : C:\HPCHEM\2\DATA\051606-F\F55863.D\ECD2B.CH
Acq On : 16 May 2006 8:07 pm Operator:
Sample : P4018 16/60 0.05 Inst : Instrumen
Misc : Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 17 9:57 2006 Quant Results File: PBQ05166.RES

Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 09:54:21 2006
Response via : Multiple Level Calibration

5-17-06



5-17-06 baseline issues were present before manual



(9) DCB (S)

10.40min 0.009ug/ml m

response 369688

(9) DCB #2 (S)

10.05min 0.009ug/ml m

response 589584

(+) = Expected Retention Time

Quantitation Report (QT Reviewed)

Signal #1 : C:\HPCHEM\2\DATA\051606-F\F55864.D\ECD1A.CH Vial: 16
Signal #2 : C:\HPCHEM\2\DATA\051606-F\F55864.D\ECD2B.CH
Acq On : 16 May 2006 8:22 pm Operator:
Sample : P4019 16/60 0.1 Inst : Instrumen
Misc : Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 17 10:04 2006 Quant Results File: PBQ05166.RES

Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 09:54:21 2006
Response via : Initial Calibration
DataAcq Meth : PSTQUICK.M

5-17-06 5-17-06

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film

Compound	RT#1	RT#2	Resp#1	Resp#2	ug/ml	ug/ml
System Monitoring Compounds						
1) S TCX	2.64	2.73	688147	815855	0.022	0.022
Spiked Amount 0.100	Range 46 - 122		Recovery =		22.00%#	22.00%#
9) S DCB	10.40	10.05	698690	1099353	0.020	0.022m
Spiked Amount 0.100	Range 26 - 86		Recovery =		20.00%#	22.00%#
Target Compounds						
2) L1 Ar1016 A	3.43	3.37	199668	227033	0.116	0.117
3) L1 Ar1016 B	3.93	3.87	361008	502332	0.111	0.117
4) L1 Ar1016 C	4.09	4.03	146603	188678	0.113	0.118
Sum Ar1016 A			707279	918043	0.340	0.352
Average Ar1016 A					0.113	0.117
5) L2 Ar1260 A	6.67	6.35	234823	325259	0.112	0.119
6) L2 Ar1260 B	7.04	6.82	274639	476109	0.109	0.121
7) L2 Ar1260 C	7.92	7.53	203415	325317	0.107	0.119
8) L2 Ar1260 D	8.49	8.59	199734	691593	0.106	0.114
Sum Ar1260 A			912612	1818277	0.433	0.473
Average Ar1260 A					0.108	0.118

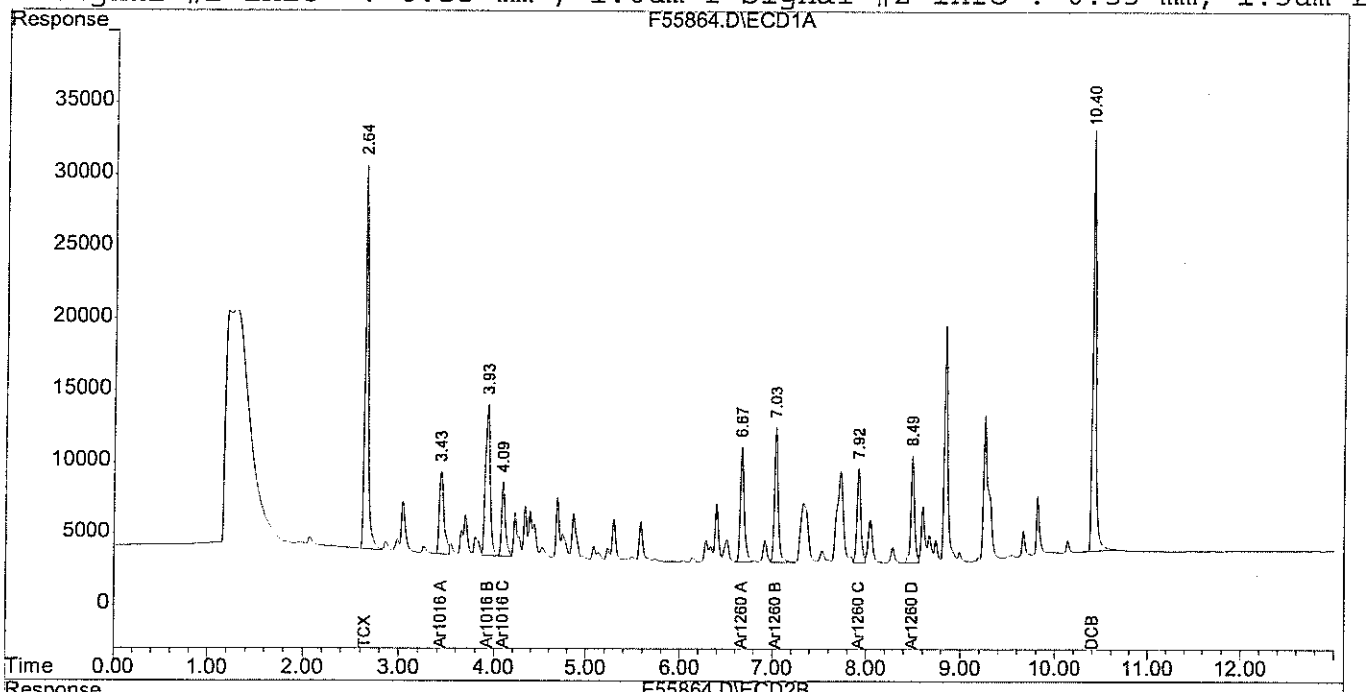
(f)=RT Delta > 1/2 Window (#)=Amounts differ by > 25% (m)=manual int.

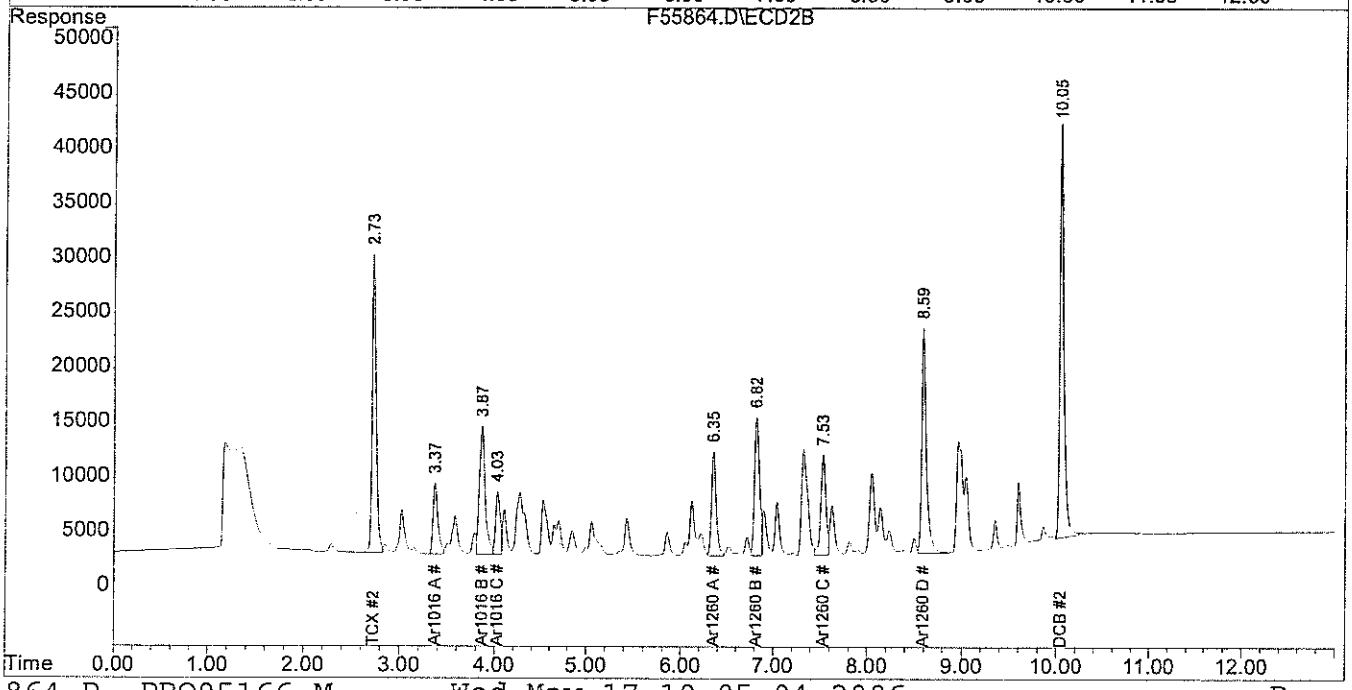
Quantitation Report (QT Reviewed)

Signal #1 : C:\HPCHEM\2\DATA\051606-F\F55864.D\ECD1A.CH Vial: 16
Signal #2 : C:\HPCHEM\2\DATA\051606-F\F55864.D\ECD2B.CH
Acq On : 16 May 2006 8:22 pm Operator:
Sample : P4019 16/60 0.1 Inst : Instrumen
Misc : Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 17 10:04 2006 Quant Results File: PBQ05166.RES

Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 09:54:21 2006
Response via : Multiple Level Calibration
DataAcq Meth : PSTQUICK.M

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film



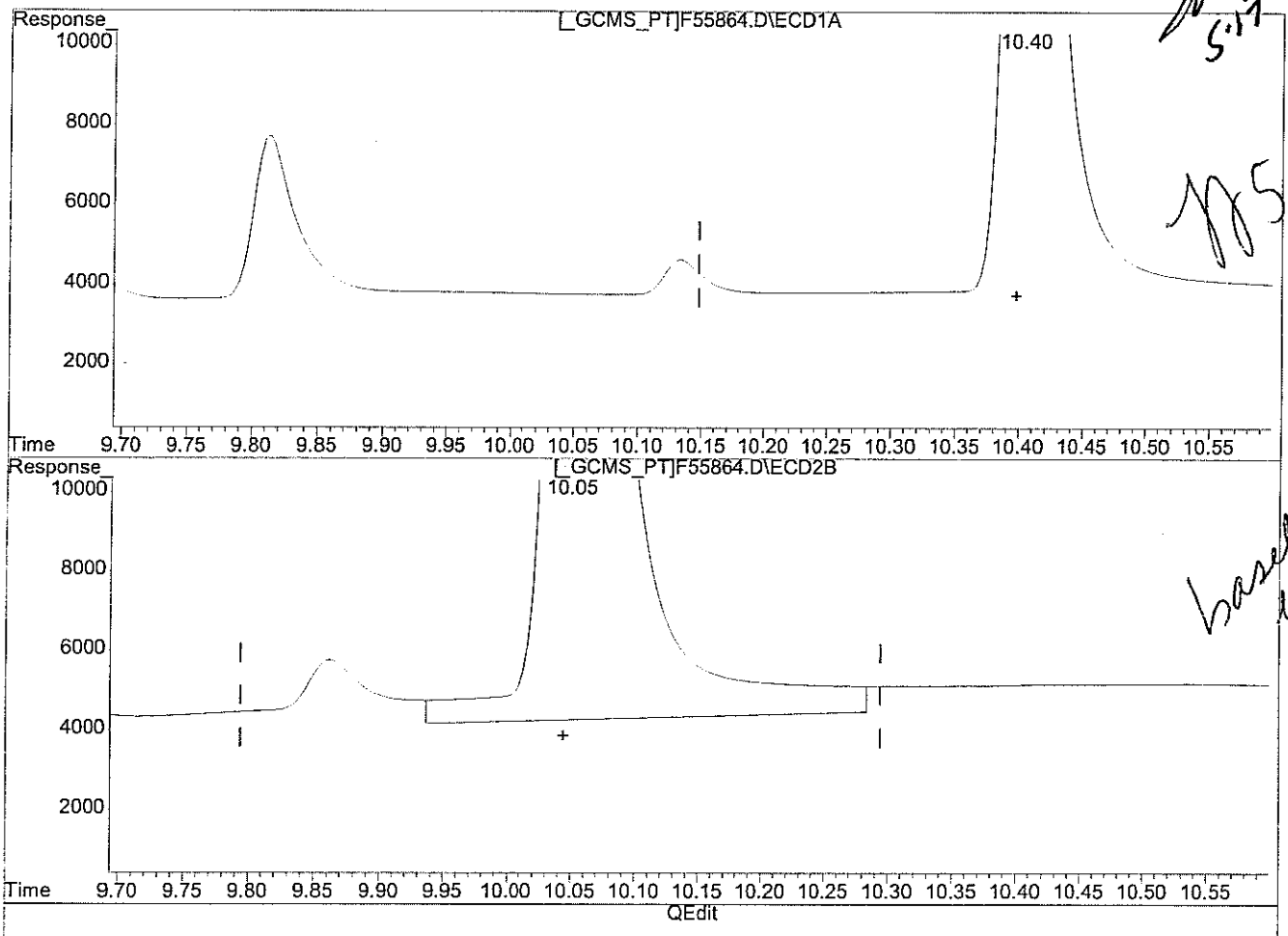


Quantitation Report (Qedit)

Signal #1 : C:\HPCHEM\2\DATA\051606-F\F55864.D\ECD1A.CH Vial: 16
Signal #2 : C:\HPCHEM\2\DATA\051606-F\F55864.D\ECD2B.CH
Acq On : 16 May 2006 8:22 pm Operator:
Sample : P4019 16/60 0.1 Inst : Instrumen
Misc : Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 17 10:04 2006 Quant Results File: PBQ05166.RES

Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 09:54:21 2006
Response via : Multiple Level Calibration

5.17.06



5.17.06

baseline/ extra area

(9) DCB (S)

10.40min 0.020ug/ml

response 698690

(9) DCB #2 (S)

10.05min 0.025ug/ml

response 1203022

(+) = Expected Retention Time

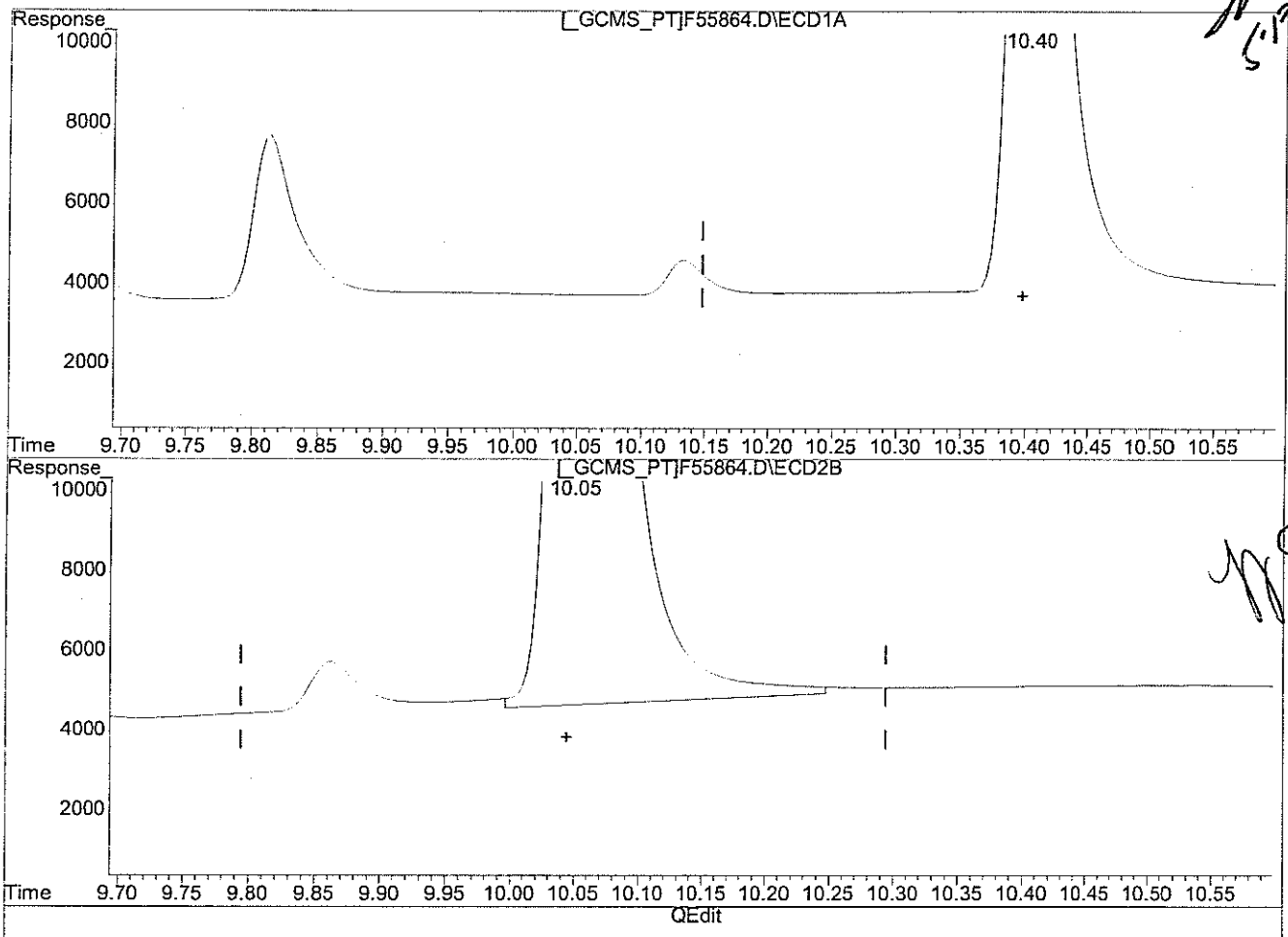
F55864.D PBQ05166.M Wed May 17 10:04:54 2006

Quantitation Report (Qedit)

Signal #1 : C:\HPCHEM\2\DATA\051606-F\F55864.D\ECD1A.CH Vial: 16
Signal #2 : C:\HPCHEM\2\DATA\051606-F\F55864.D\ECD2B.CH
Acq On : 16 May 2006 8:22 pm Operator:
Sample : P4019 16/60 0.1 Inst : Instrumen
Misc : Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 17 10:04 2006 Quant Results File: PBQ05166.RES

Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 09:54:21 2006
Response via : Multiple Level Calibration

5.19.06



5.17.06

(9) DCB (S)

10.40min 0.020ug/ml

response 698690

(9) DCB #2 (S)

10.05min 0.022ug/ml m

response 1099353

(+) = Expected Retention Time

Quantitation Report (QT Reviewed)

Signal #1 : C:\HPCHEM\2\DATA\051606-F\F55865.D\ECD1A.CH Vial: 17
Signal #2 : C:\HPCHEM\2\DATA\051606-F\F55865.D\ECD2B.CH
Acq On : 16 May 2006 8:36 pm Operator:
Sample : P4020 16/60 0.25 Inst : Instrumen
Misc : Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 17 9:54 2006 Quant Results File: PBQ05166.RES

Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 09:54:21 2006
Response via : Initial Calibration
DataAcq Meth : PSTQUICK.M

5-17-06 5-17-06

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film

Compound	RT#1	RT#2	Resp#1	Resp#2	ug/ml	ug/ml
System Monitoring Compounds						
1) S TCX	2.64	2.73	1837346	2057065	0.058	0.058
Spiked Amount 0.100	Range 46 - 122			Recovery =	58.00%	58.00%
9) S DCB	10.40	10.04	1671904	2550743	0.054	0.061
Spiked Amount 0.100	Range 26 - 86			Recovery =	54.00%	61.00%
Target Compounds						
2) L1 Ar1016 A	3.44	3.37	469862	514274	0.296	0.300
3) L1 Ar1016 B	3.93	3.87	885341	1150893	0.293	0.299
4) L1 Ar1016 C	4.09	4.03	365795	438601	0.294	0.304
Sum Ar1016 A			1720998	2103768	0.884	0.903
Average Ar1016 A					0.295	0.301
5) L2 Ar1260 A	6.67	6.35	562785	744547	0.288	0.321
6) L2 Ar1260 B	7.03	6.82	670680	1091231	0.285	0.319
7) L2 Ar1260 C	7.92	7.53	510810	774876	0.279	0.318
8) L2 Ar1260 D	8.49	8.59	500733	1688943	0.278	0.299
Sum Ar1260 A			2245008	4299597	1.130	1.256
Average Ar1260 A					0.283	0.314

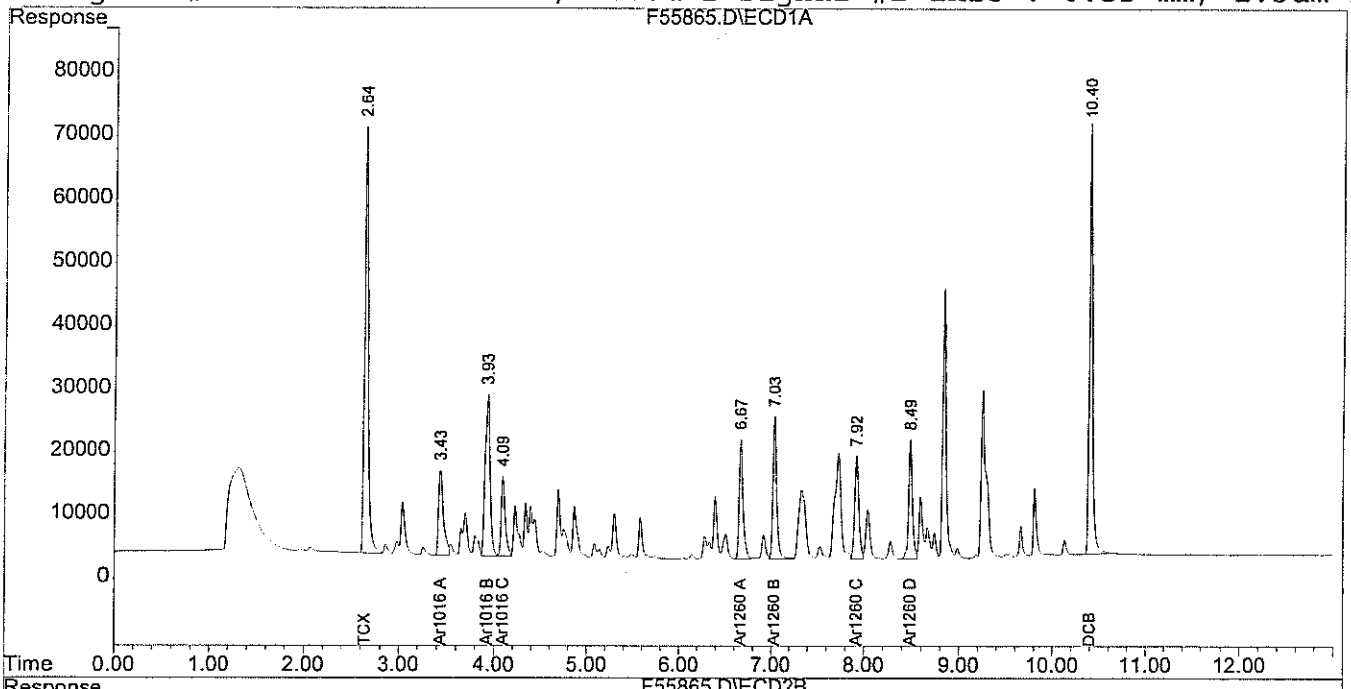
(f)=RT Delta > 1/2 Window (#)=Amounts differ by > 25% (m)=manual int.

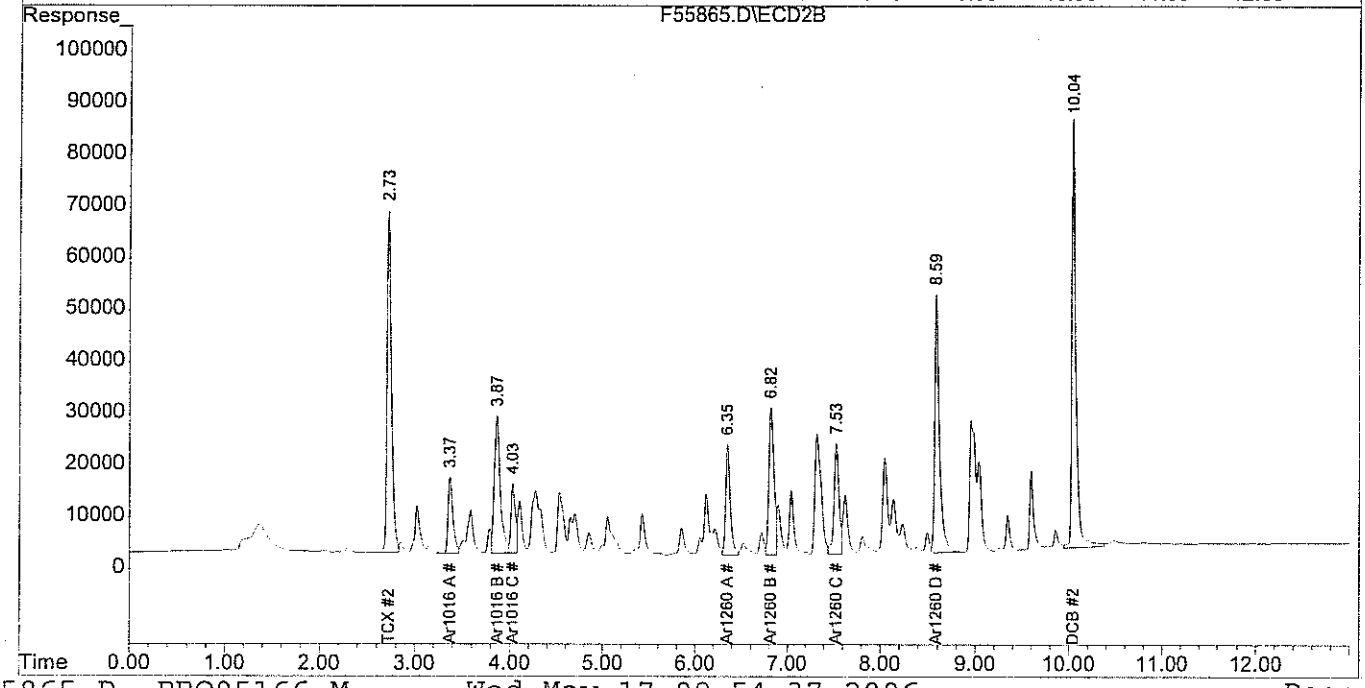
Quantitation Report (QT Reviewed)

Signal #1 : C:\HPCHEM\2\DATA\051606-F\F55865.D\ECD1A.CH Vial: 17
Signal #2 : C:\HPCHEM\2\DATA\051606-F\F55865.D\ECD2B.CH
Acq On : 16 May 2006 8:36 pm Operator:
Sample : P4020 16/60 0.25 Inst : Instrumen
Misc : Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 17 9:54 2006 Quant Results File: PBQ05166.RES

Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 09:54:21 2006
Response via : Multiple Level Calibration
DataAcq Meth : PSTQUICK.M

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film





Quantitation Report (QT Reviewed)

Signal #1 : C:\HPCHEM\2\DATA\051606-F\F55866.D\ECD1A.CH Vial: 18
Signal #2 : C:\HPCHEM\2\DATA\051606-F\F55866.D\ECD2B.CH
Acq On : 16 May 2006 8:51 pm Operator:
Sample : P4021 16/60 0.5 Inst : Instrumen
Misc : Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 17 10:05 2006 Quant Results File: PBQ05166.RES

Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 09:54:21 2006
Response via : Initial Calibration
DataAcq Meth : PSTQUICK.M

5.17.06 5.17.06

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film

Compound	RT#1	RT#2	Resp#1	Resp#2	ug/ml	ug/ml
System Monitoring Compounds						
1) S TCX	2.64	2.73	3525739	3853418	0.112	0.110
Spiked Amount 0.100	Range 46 - 122		Recovery =		112.00%	110.00%
9) S DCB	10.39	10.04	3091651	4358733	0.104	0.108
Spiked Amount 0.100	Range 26 - 86		Recovery =		104.00%#	108.00%#
Target Compounds						
2) L1 Ar1016 A	3.43	3.37	867015	920329	0.561	0.558
L1 Ar1016 B	3.93	3.87	1650244	2072594	0.559	0.557
4) L1 Ar1016 C	4.09	4.03	686360	787671	0.560	0.564
Sum Ar1016 A			3203619	3780595	1.680	1.679
Average Ar1016 A					0.560	0.560
5) L2 Ar1260 A	6.67	6.35	1071560	1271657	0.562	0.575
6) L2 Ar1260 B	7.03	6.82	1268608	1900394	0.552	0.579
7) L2 Ar1260 C	7.92	7.53	993398	1384156	0.549	0.587
8) L2 Ar1260 D	8.49	8.59	972866	3155830	0.547	0.571
Sum Ar1260 A			4306432	7712037	2.210	2.312
Average Ar1260 A					0.553	0.578

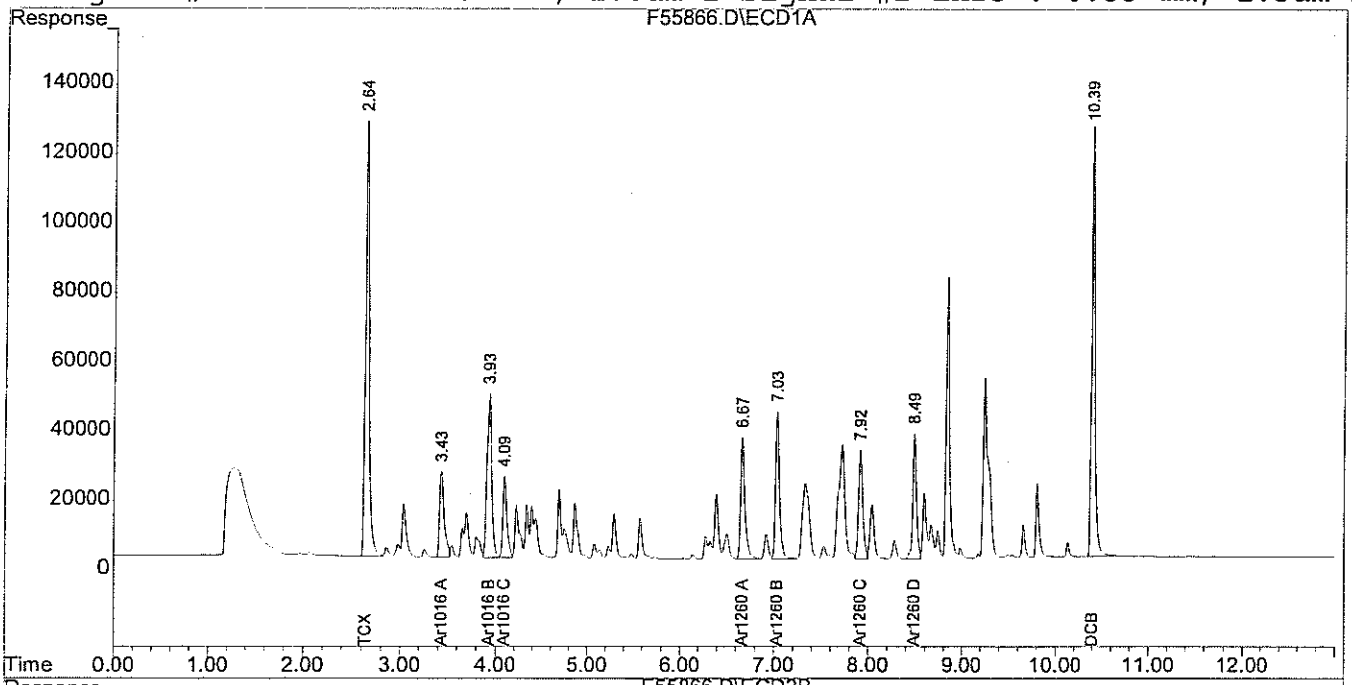
(f)=RT Delta > 1/2 Window (#)=Amounts differ by > 25% (m)=manual int

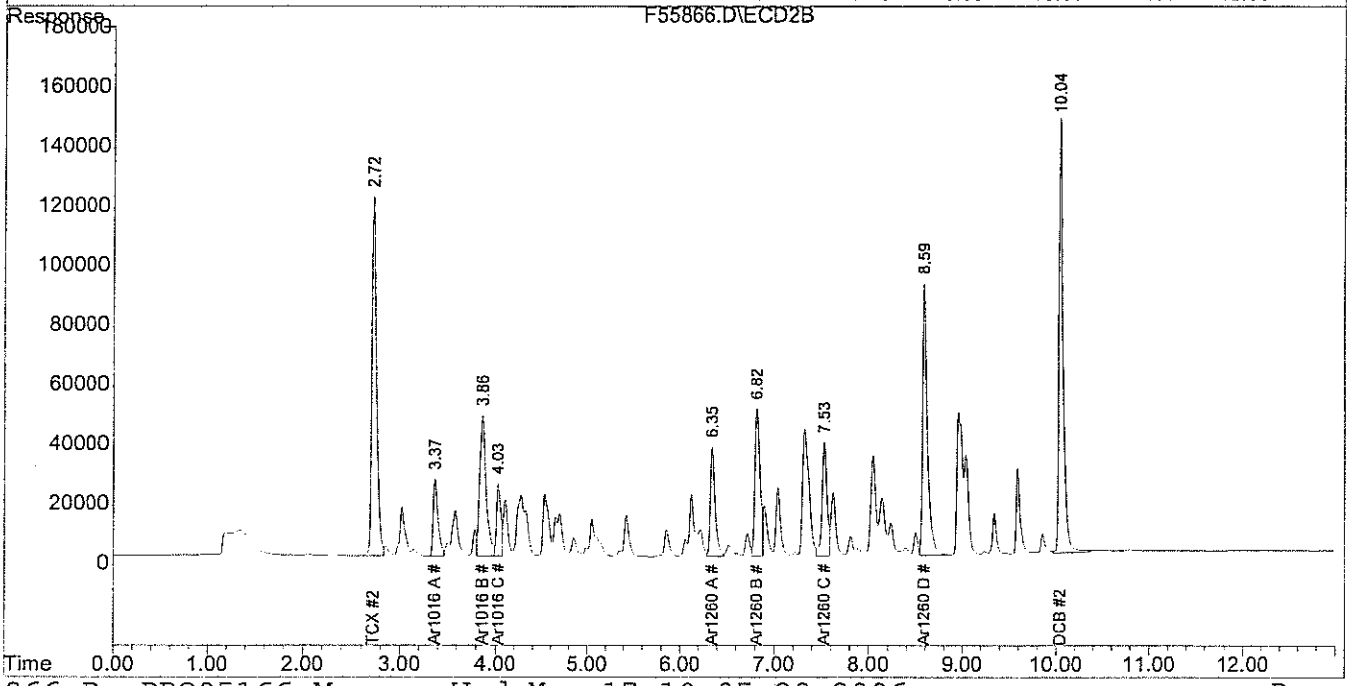
Quantitation Report (QT Reviewed)

Signal #1 : C:\HPCHEM\2\DATA\051606-F\F55866.D\ECD1A.CH Vial: 18
Signal #2 : C:\HPCHEM\2\DATA\051606-F\F55866.D\ECD2B.CH
Acq On : 16 May 2006 8:51 pm Operator:
Sample : P4021 16/60 0.5 Inst : Instrumen
Misc : Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 17 10:05 2006 Quant Results File: PBQ05166.RES

Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 09:54:21 2006
Response via : Multiple Level Calibration
DataAcq Meth : PSTQUICK.M

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film





Quantitation Report (QT Reviewed)

Signal #1 : C:\HPCHEM\2\DATA\051606-F\F55867.D\ECD1A.CH Vial: 19
Signal #2 : C:\HPCHEM\2\DATA\051606-F\F55867.D\ECD2B.CH
Acq On : 16 May 2006 9:20 pm Operator:
Sample : P4022 16/60 1.0 Inst : Instrumen
Misc : Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 17 10:05 2006 Quant Results File: PBQ05166.RES

Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 09:54:21 2006
Response via : Initial Calibration
DataAcq Meth : PSTQUICK.M

5.17.06 5.17.06

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film

Compound	RT#1	RT#2	Resp#1	Resp#2	ug/ml	ug/ml
System Monitoring Compounds						
1) S TCX	2.64	2.73	6920671	7382493	0.219	0.211
Spiked Amount 0.100	Range 46 - 122		Recovery =		219.00%#	211.00%#
9) S DCB	10.40	10.04	5840708	7809116	0.200	0.200
Spiked Amount 0.100	Range 26 - 86		Recovery =		200.00%#	200.00%#
Target Compounds						
2) L1 Ar1016 A	3.43	3.37	1663419	1708138	1.092	1.059
3) L1 Ar1016 B	3.93	3.87	3193003	3862796	1.095	1.058
4) L1 Ar1016 C	4.09	4.03	1338659	1462474	1.101	1.068
Sum Ar1016 A			6195082	7033407	3.288	3.185
Average Ar1016 A					1.096	1.062
5) L2 Ar1260 A	6.67	6.35	2047135	2313020	1.087	1.076
6) L2 Ar1260 B	7.03	6.82	2456273	3457427	1.083	1.080
7) L2 Ar1260 C	7.92	7.53	1964211	2559137	1.092	1.107
8) L2 Ar1260 D	8.49	8.59	1930282	5977809	1.093	1.094
Sum Ar1260 A			8397902	14307393	4.355	4.357
Average Ar1260 A					1.089	1.089

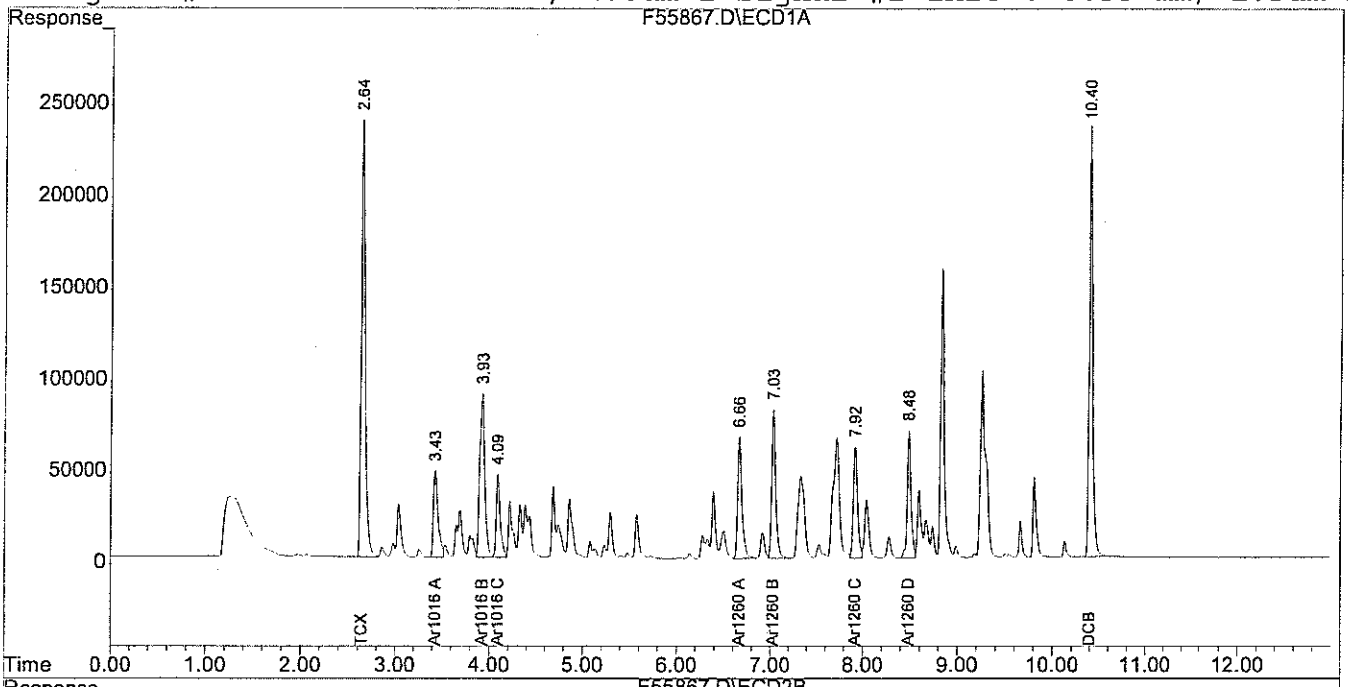
(f)=RT Delta > 1/2 Window (#)=Amounts differ by > 25% (m)=manual int.

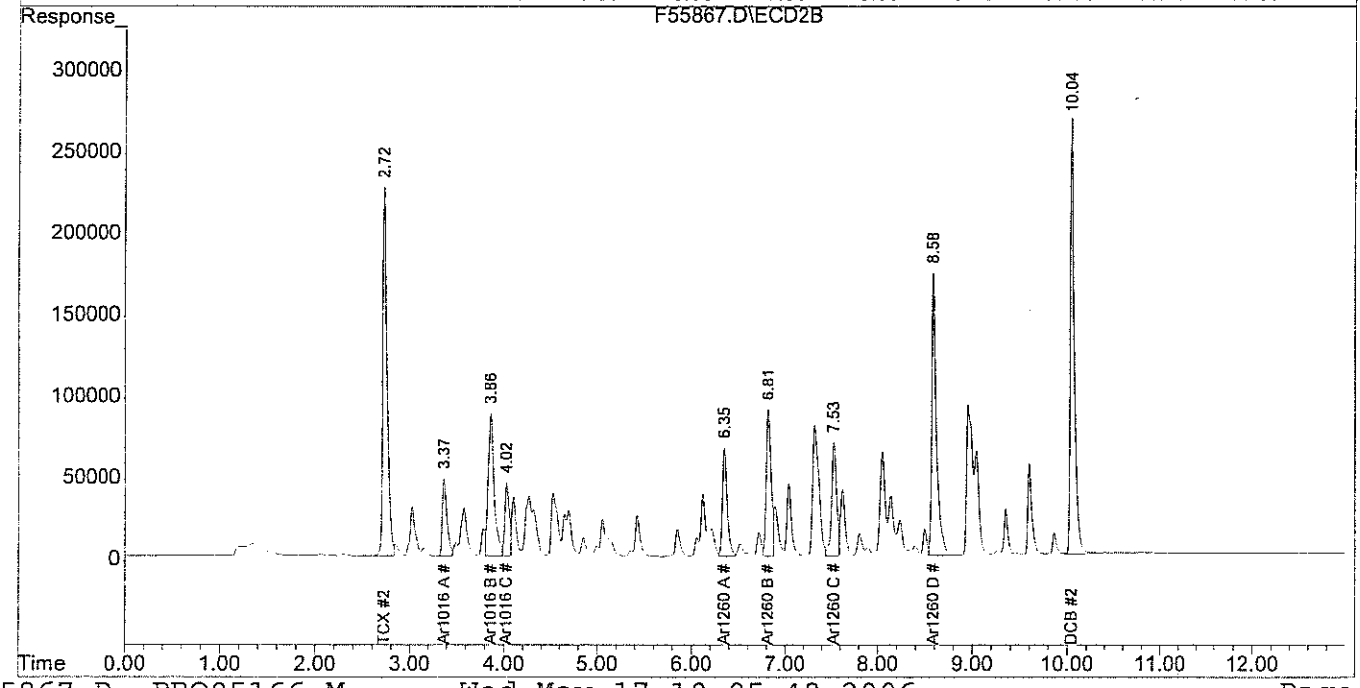
Quantitation Report (QT Reviewed)

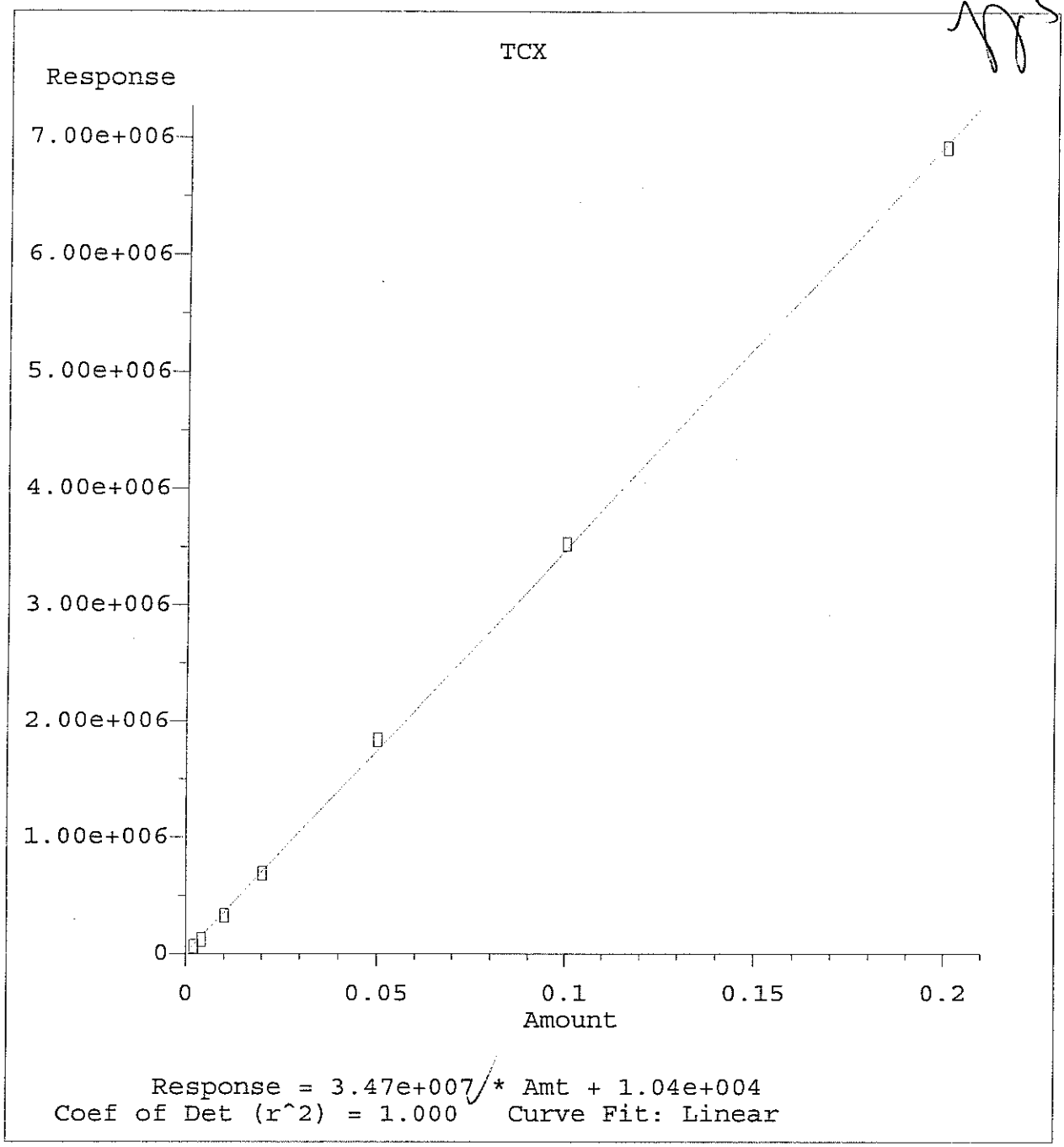
Signal #1 : C:\HPCHEM\2\DATA\051606-F\F55867.D\ECD1A.CH Vial: 19
Signal #2 : C:\HPCHEM\2\DATA\051606-F\F55867.D\ECD2B.CH
Acq On : 16 May 2006 9:20 pm Operator:
Sample : P4022 16/60 1.0 Inst : Instrumen
Misc : Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 17 10:05 2006 Quant Results File: PBQ05166.RES

Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 09:54:21 2006
Response via : Multiple Level Calibration
DataAcq Meth : PSTQUICK.M

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film

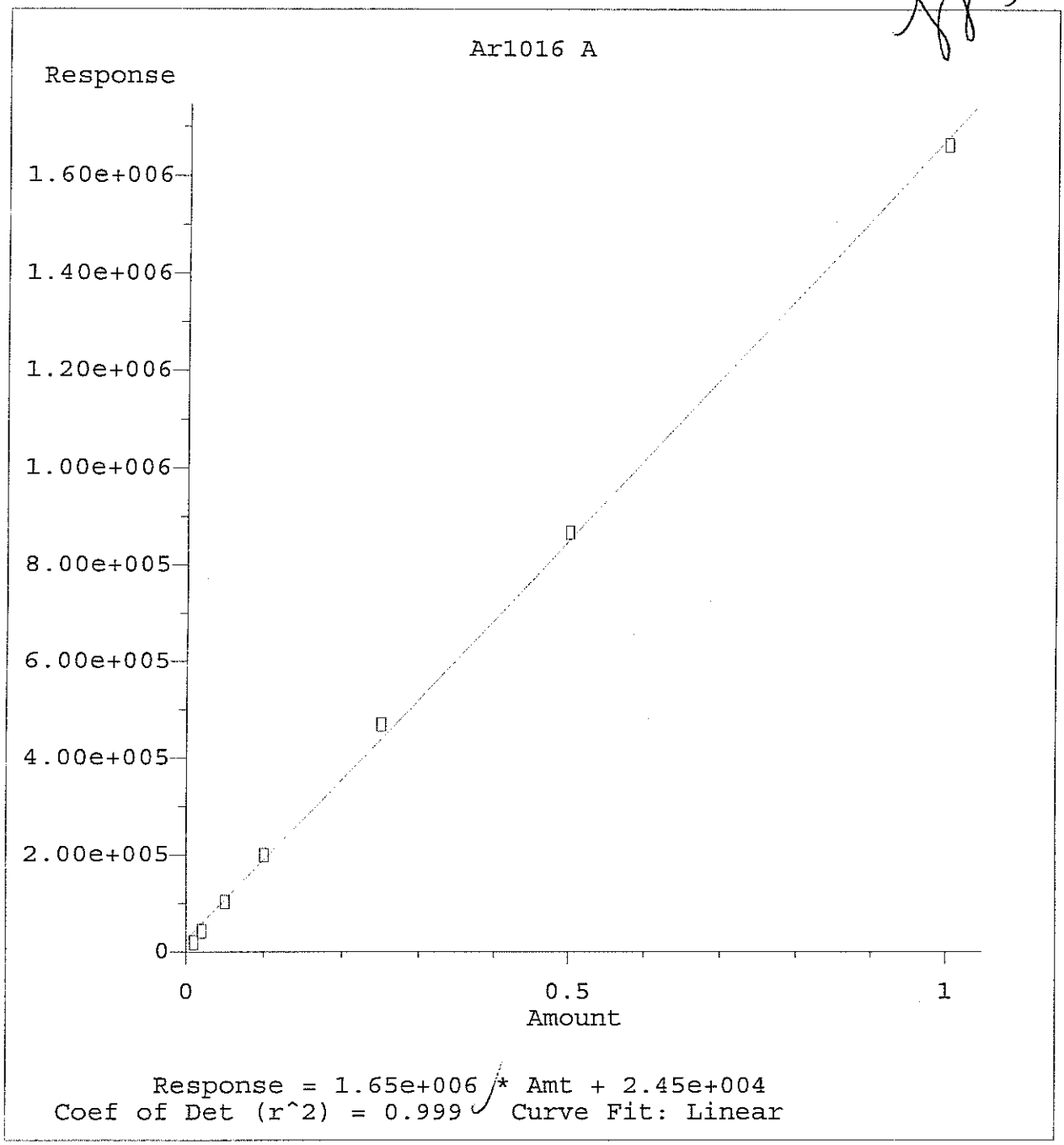






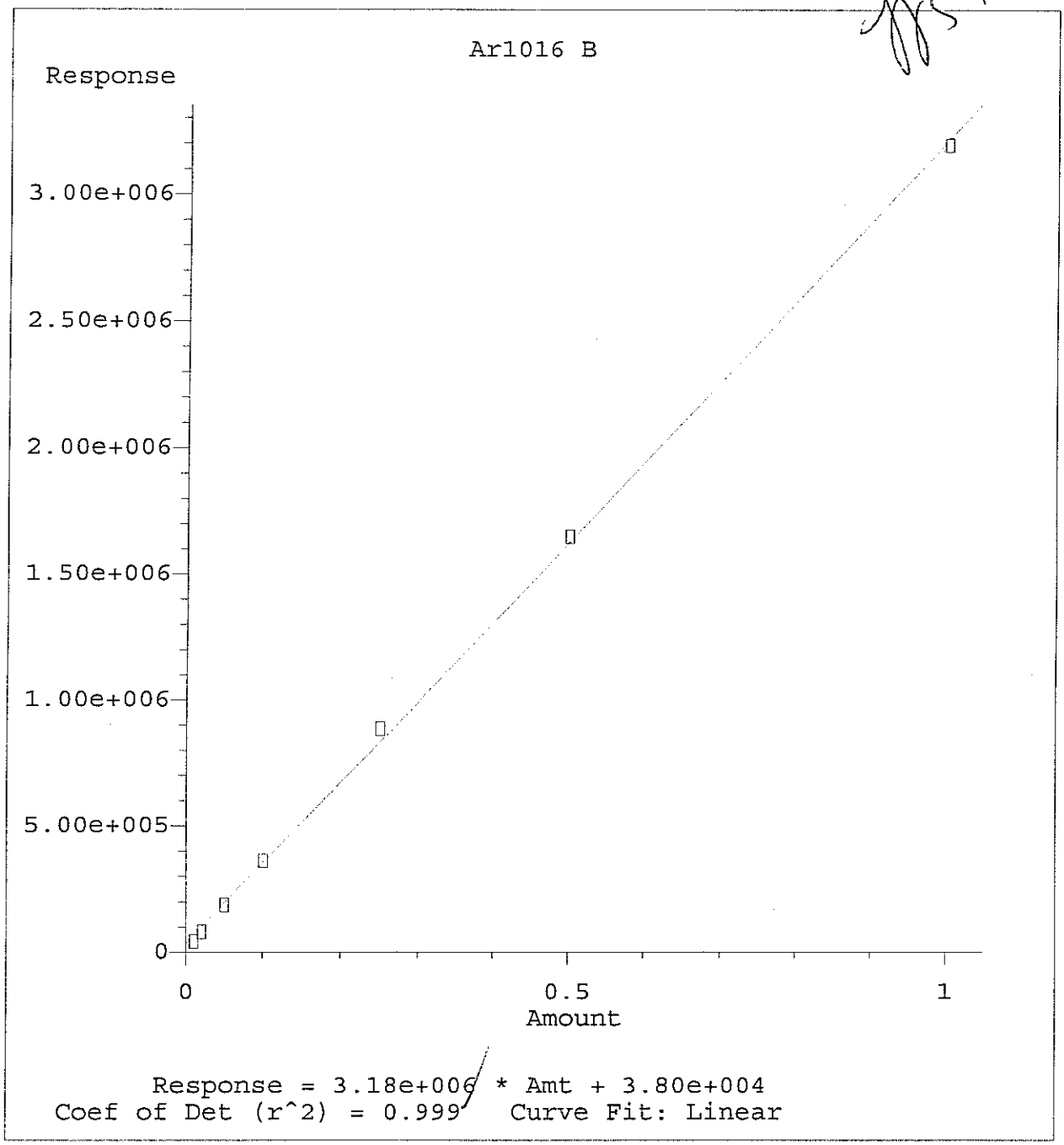
Method Name: C:\HPCHEM\2\METHODS\PBQ05166.M
Calibration Table Last Updated: Wed May 17 10:06:32 2006

5.17.06



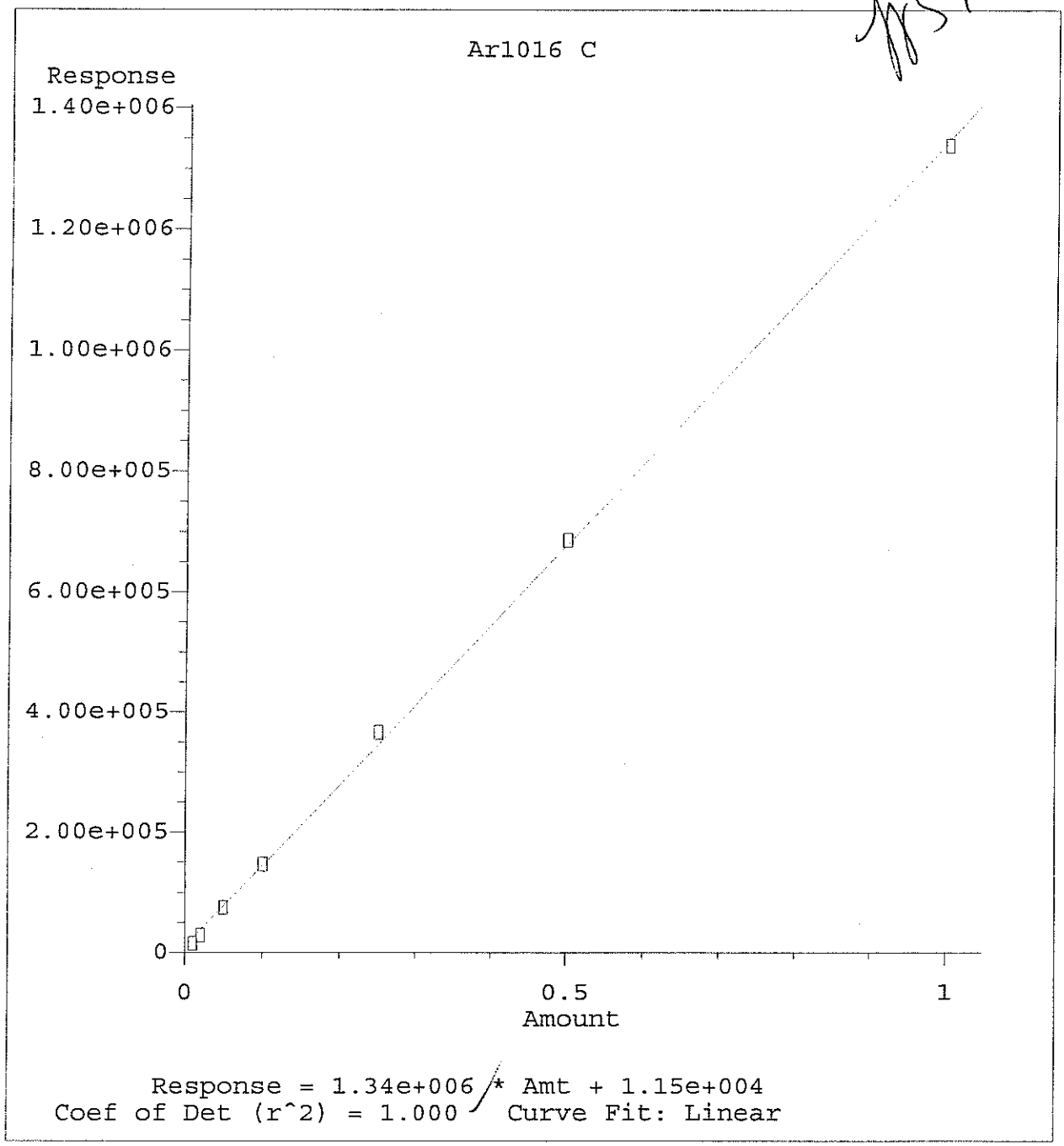
Method Name: C:\HPCHEM\2\METHODS\PBQ05166.M
Calibration Table Last Updated: Wed May 17 10:06:32 2006

5.17.06



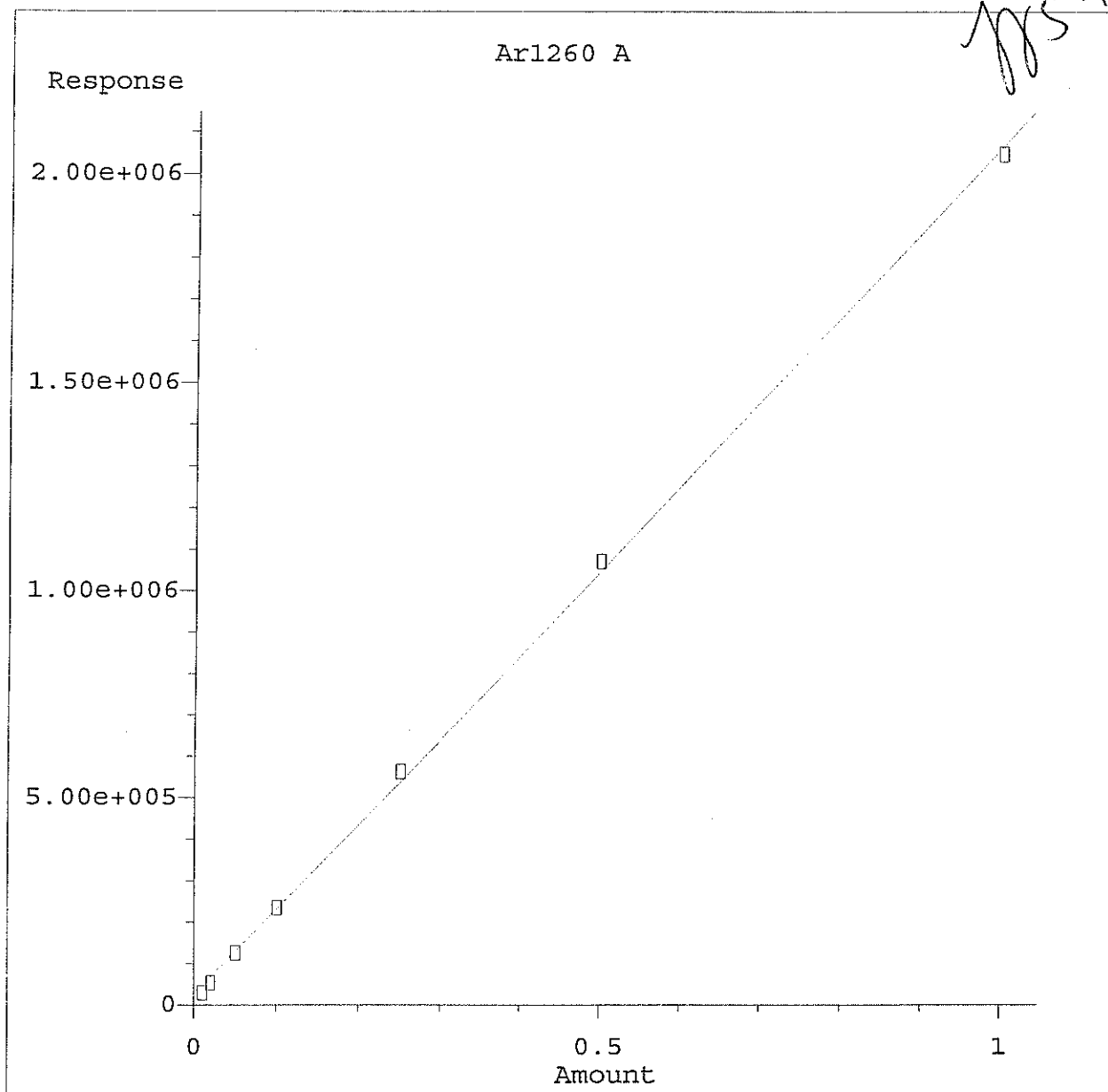
Method Name: C:\HPCHEM\2\METHODS\PBQ05166.M
Calibration Table Last Updated: Wed May 17 10:06:32 2006

5.17.06



Method Name: C:\HPCHEM\2\METHODS\PBQ05166.M
Calibration Table Last Updated: Wed May 17 10:06:32 2006

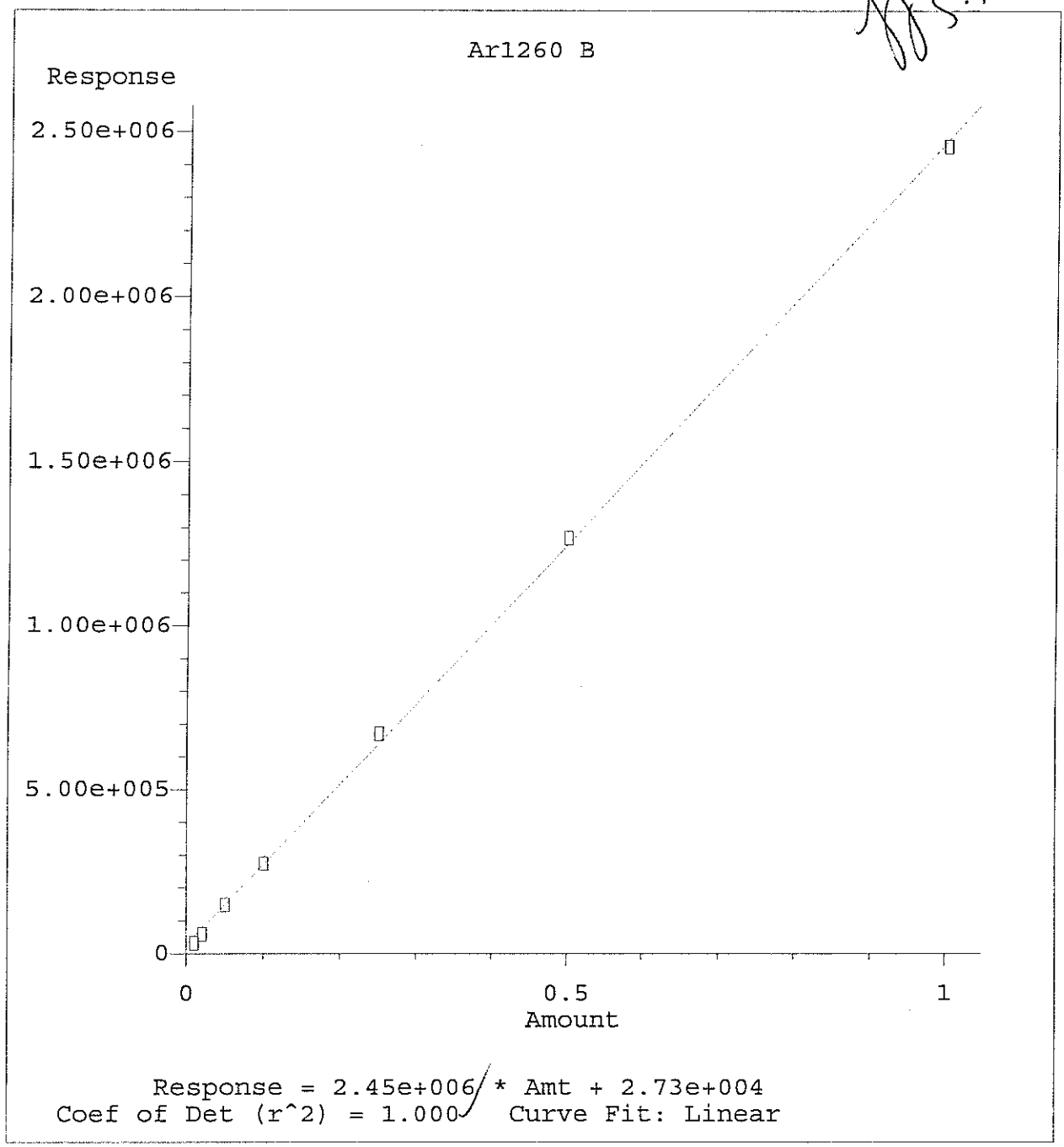
JJS 5.17.06



Response = 2.04e+006 * Amt + 2.75e+004
Coef of Det (r^2) = 0.999 Curve Fit: Linear

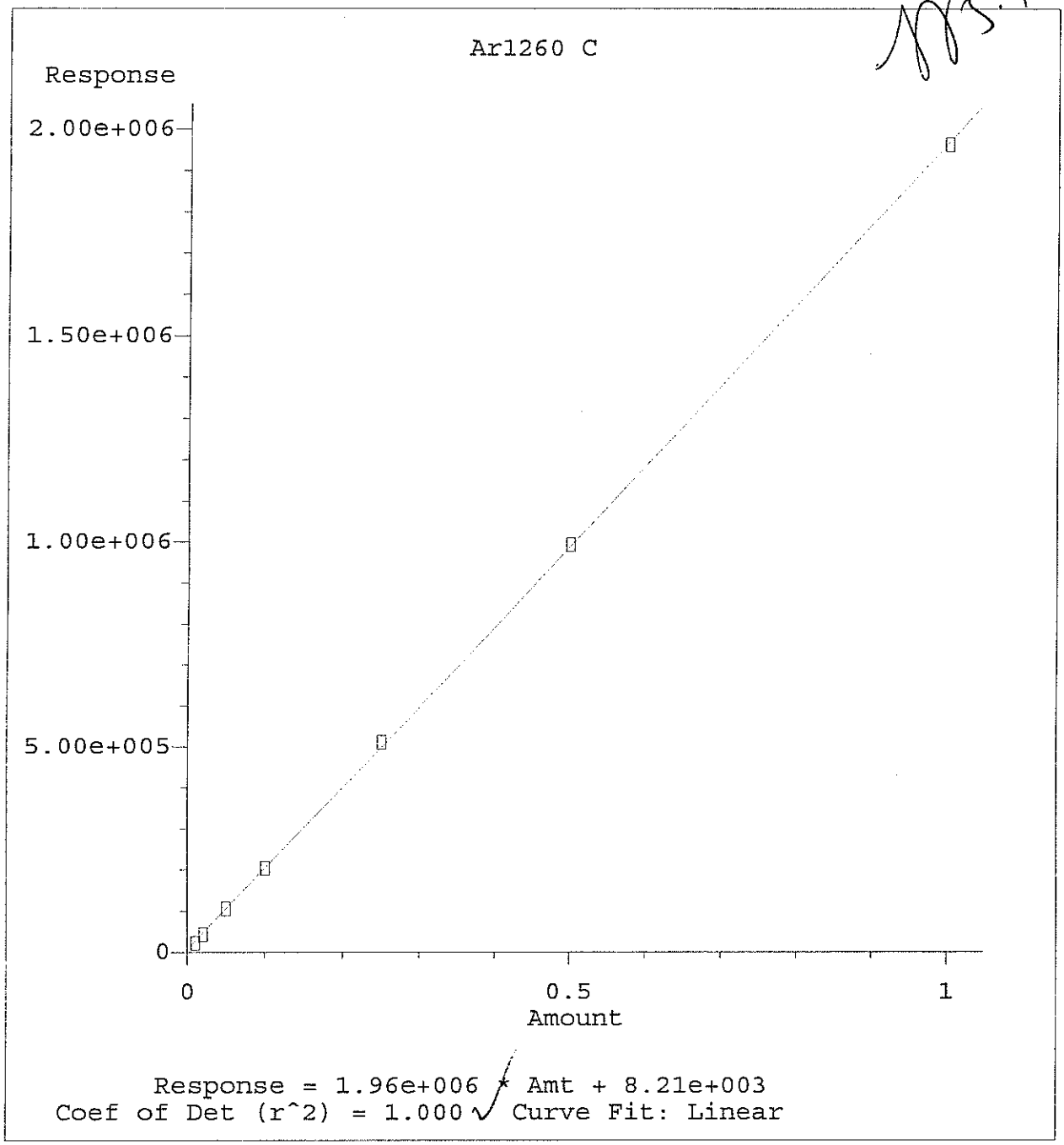
Method Name: C:\HPCHEM\2\METHODS\PBQ05166.M
Calibration Table Last Updated: Wed May 17 10:06:32 2006

5.17.06



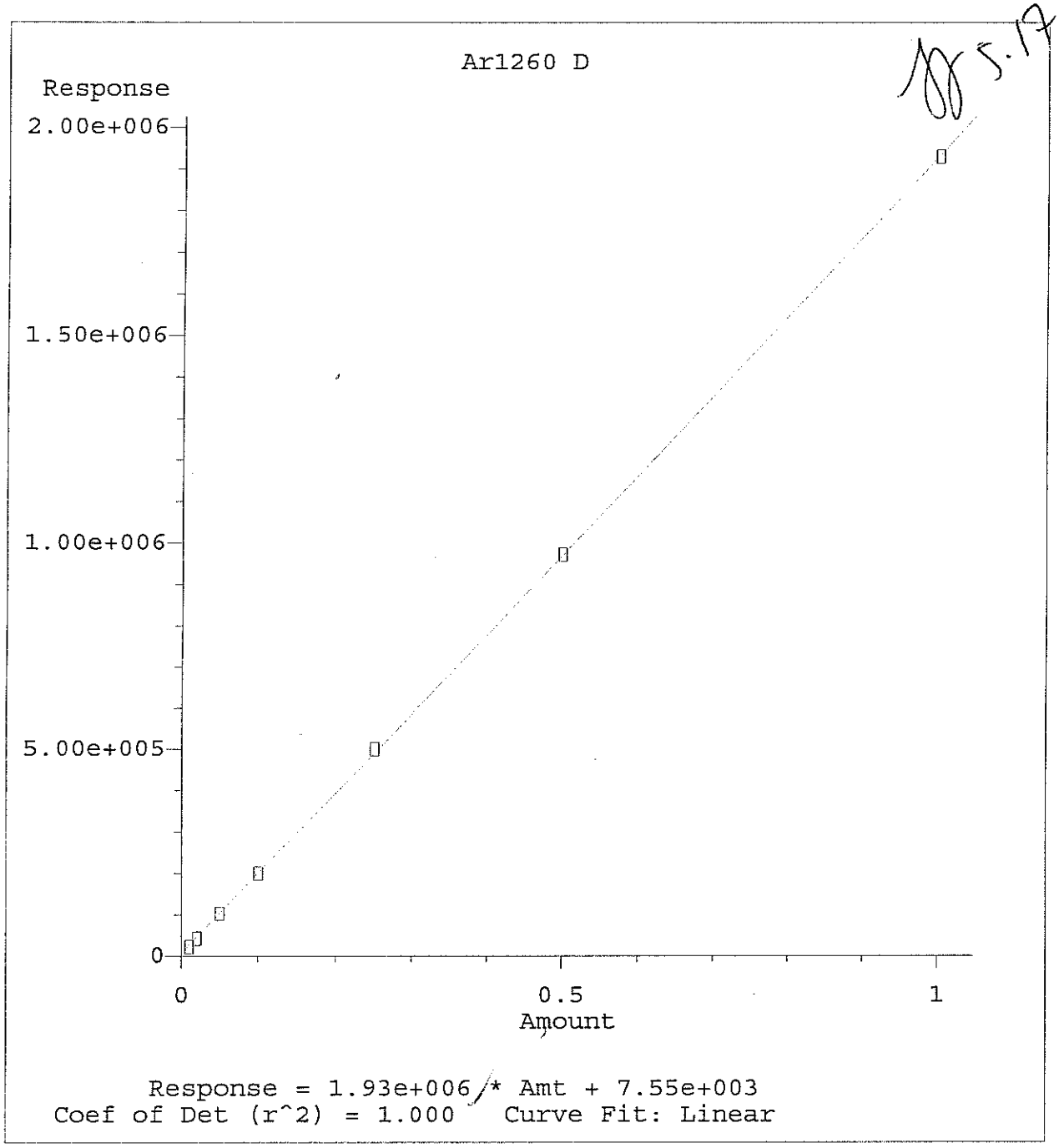
Method Name: C:\HPCHEM\2\METHODS\PBQ05166.M
Calibration Table Last Updated: Wed May 17 10:06:32 2006

5.17.06



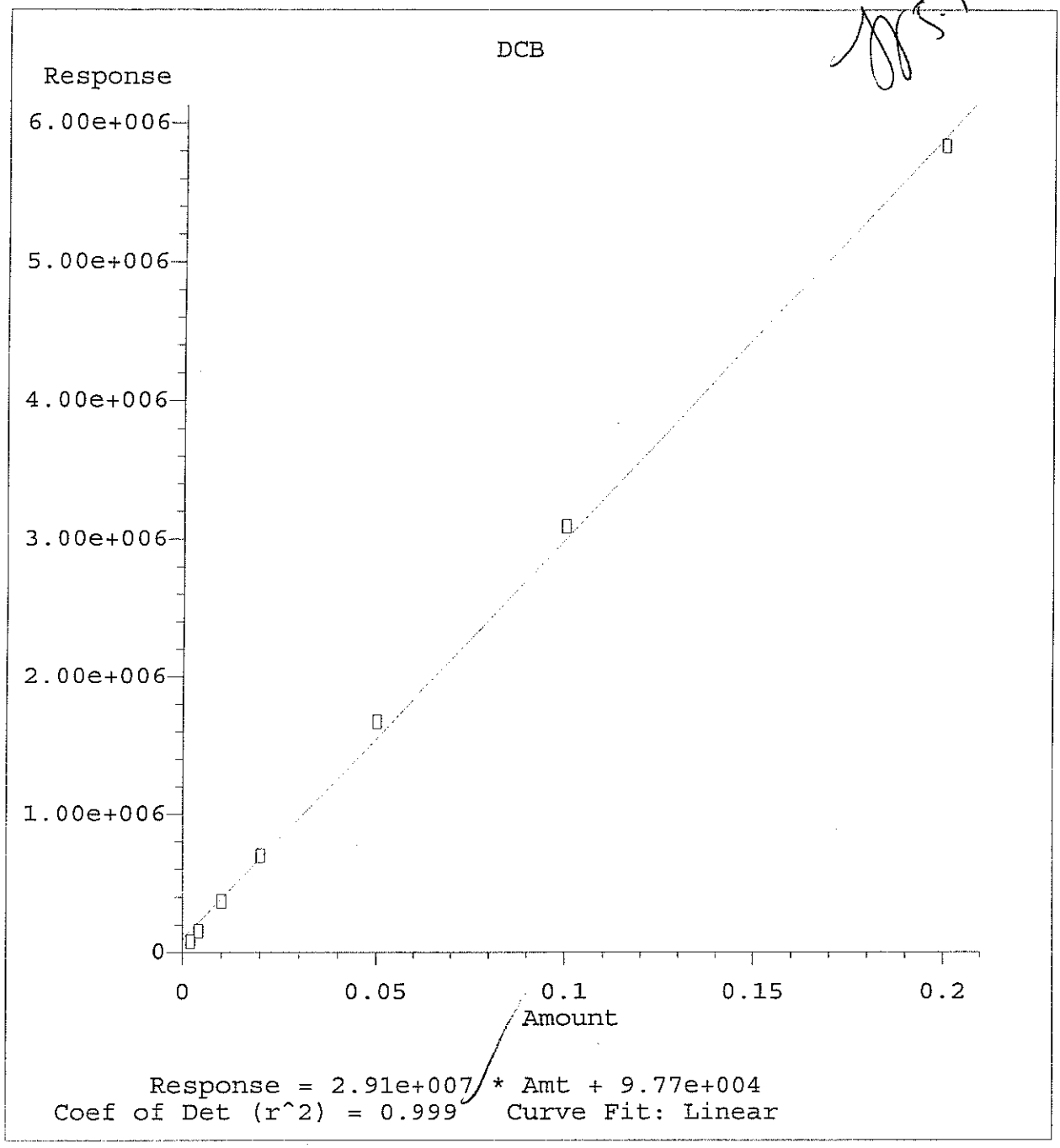
Method Name: C:\HPCHEM\2\METHODS\PBQ05166.M
Calibration Table Last Updated: Wed May 17 10:06:32 2006

5.17.06

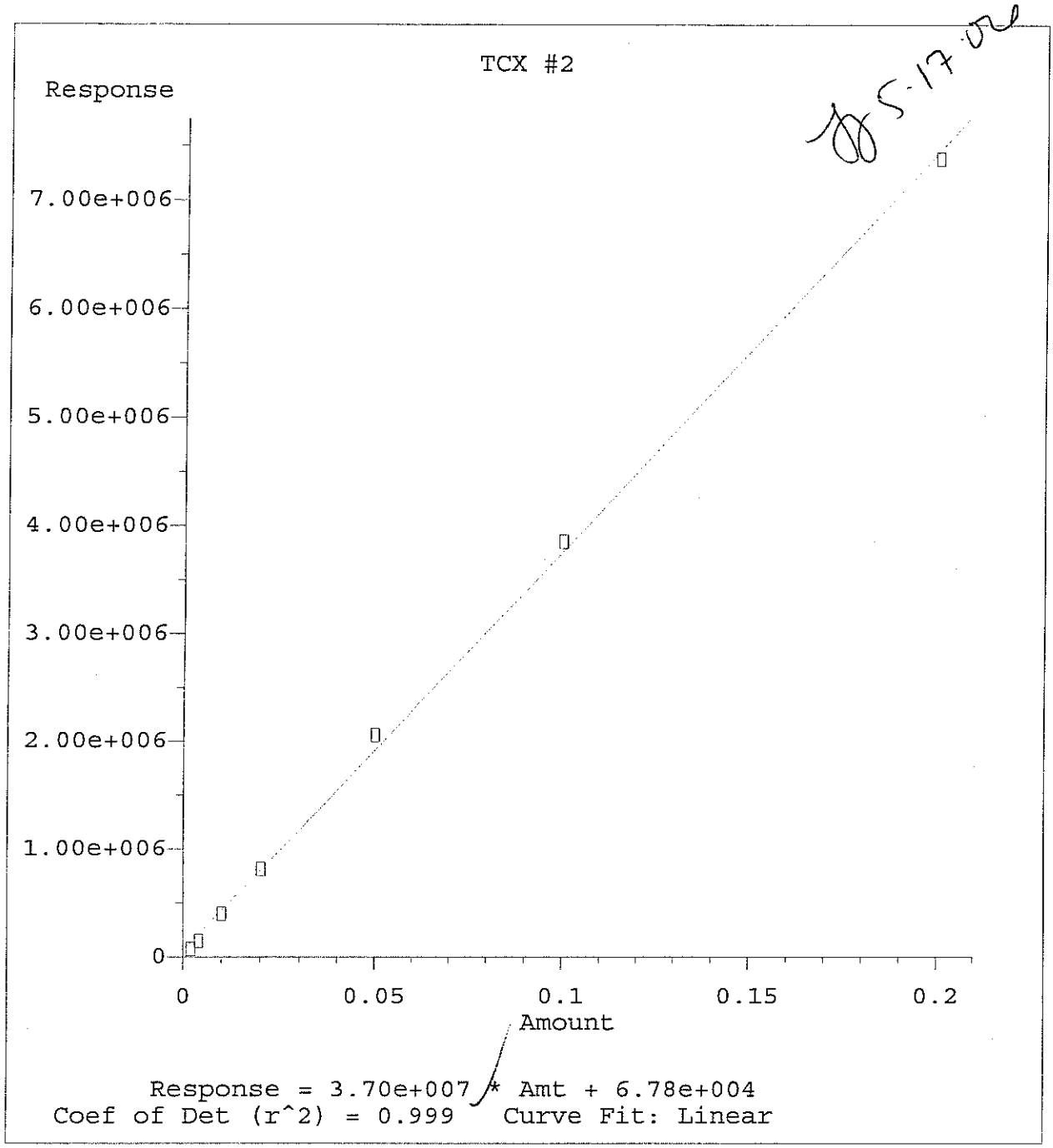


Method Name: C:\HPCHEM\2\METHODS\PBQ05166.M
Calibration Table Last Updated: Wed May 17 10:06:32 2006

5-17-06

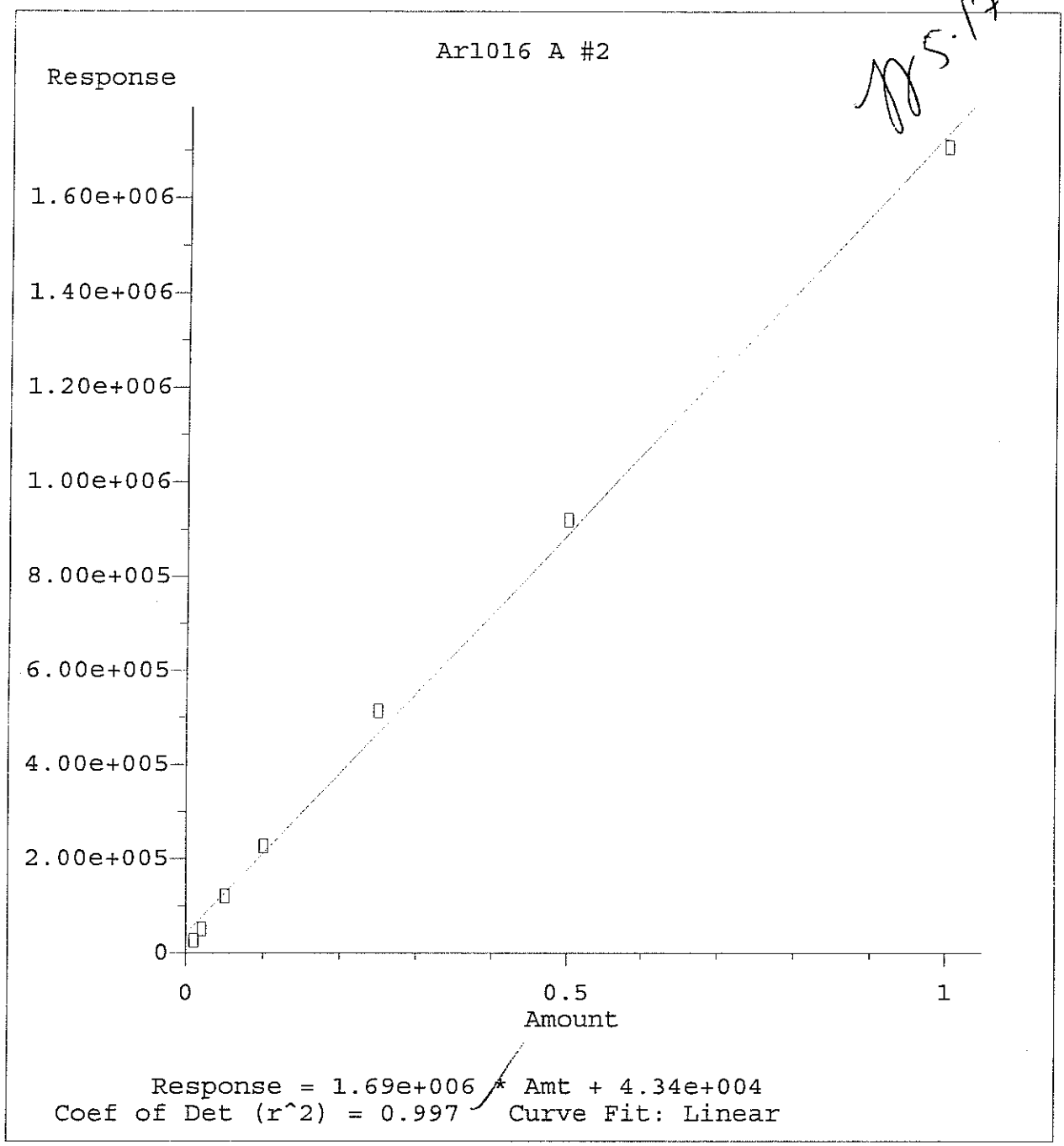


Method Name: C:\HPCHEM\2\METHODS\PBQ05166.M
Calibration Table Last Updated: Wed May 17 10:06:32 2006

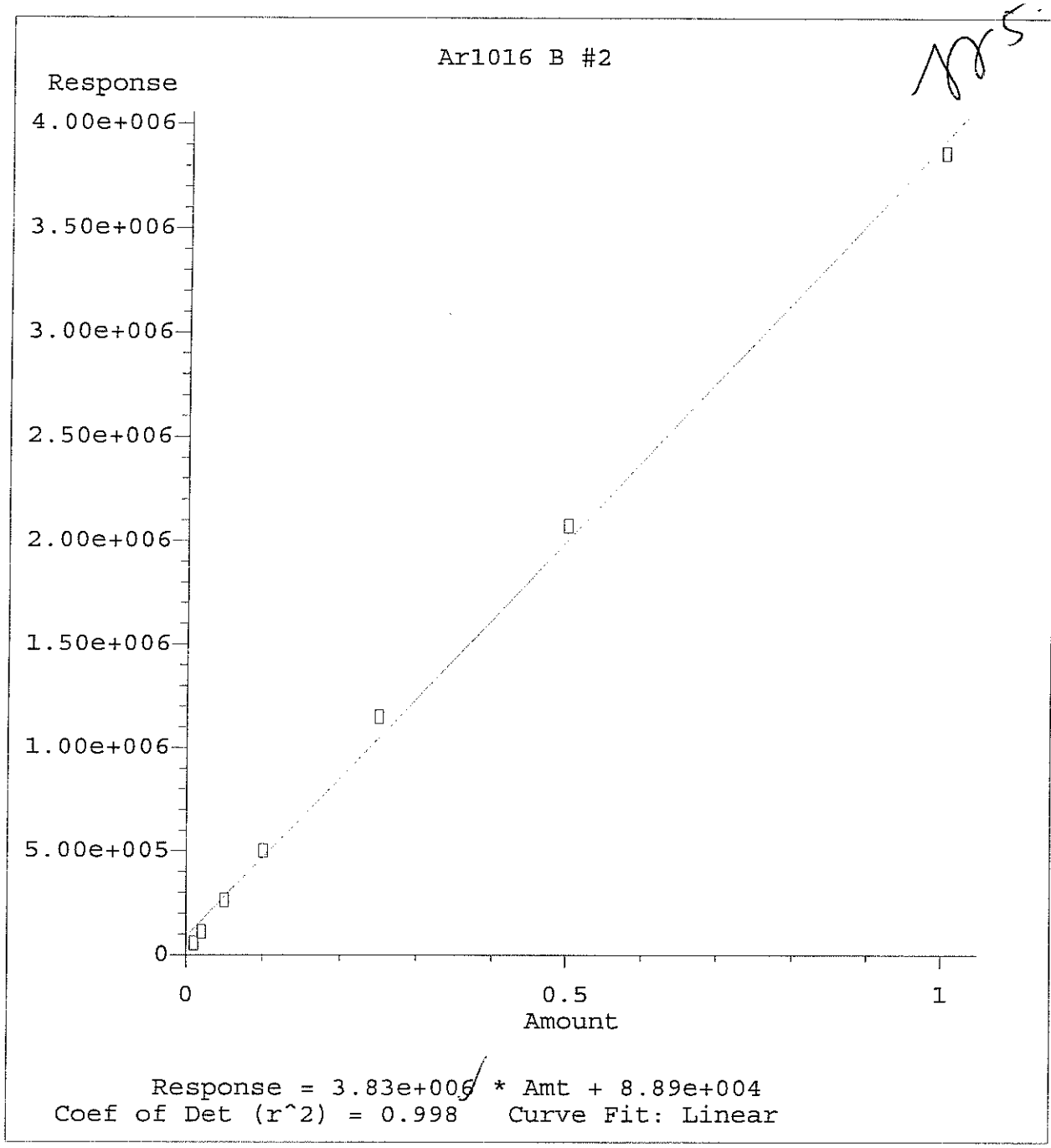


Method Name: C:\HPCHEM\2\METHODS\PBQ05166.M
Calibration Table Last Updated: Wed May 17 10:06:32 2006

5.17.06

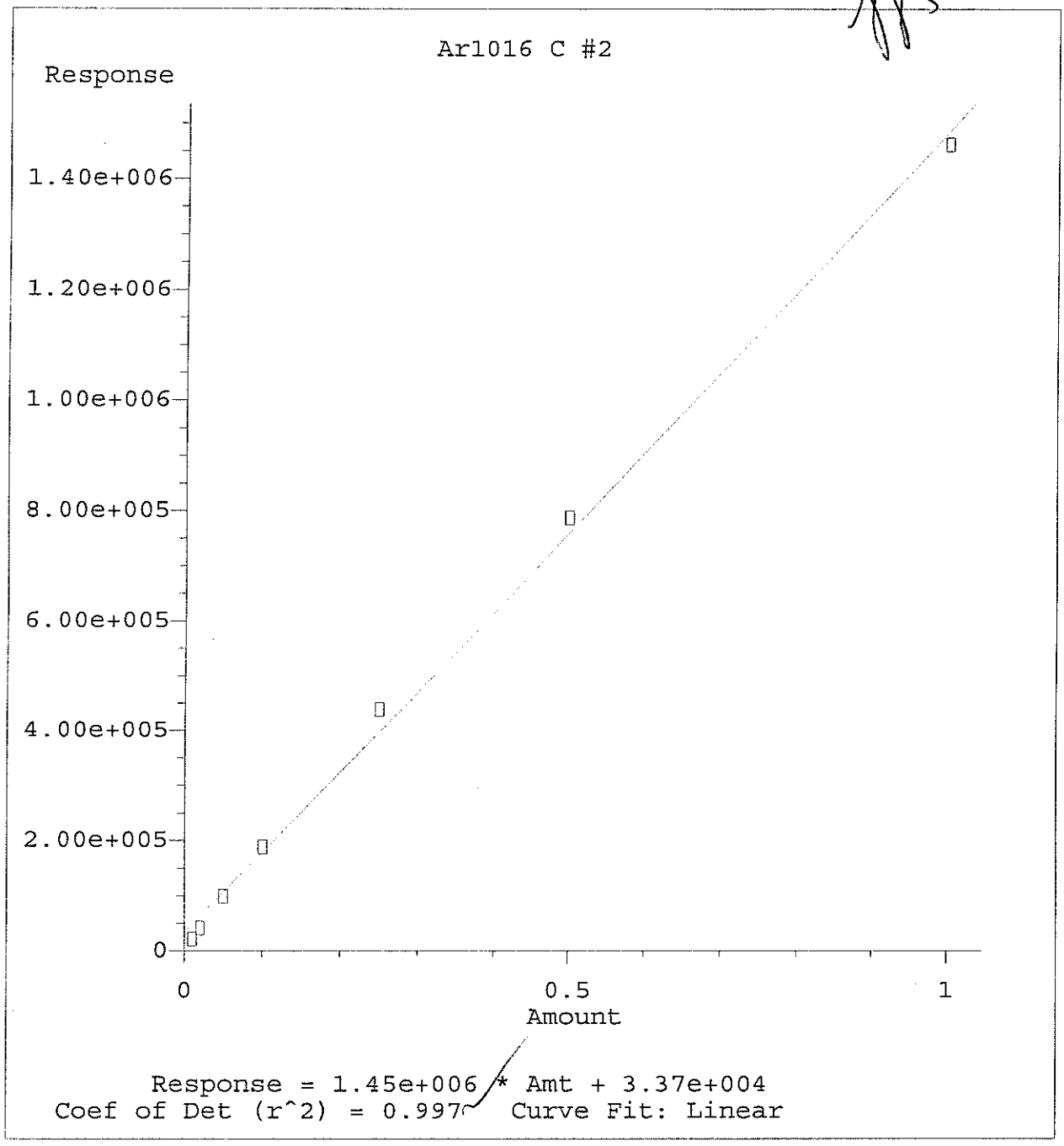


Method Name: C:\HPCHEM\2\METHODS\PBQ05166.M
Calibration Table Last Updated: Wed May 17 10:06:32 2006



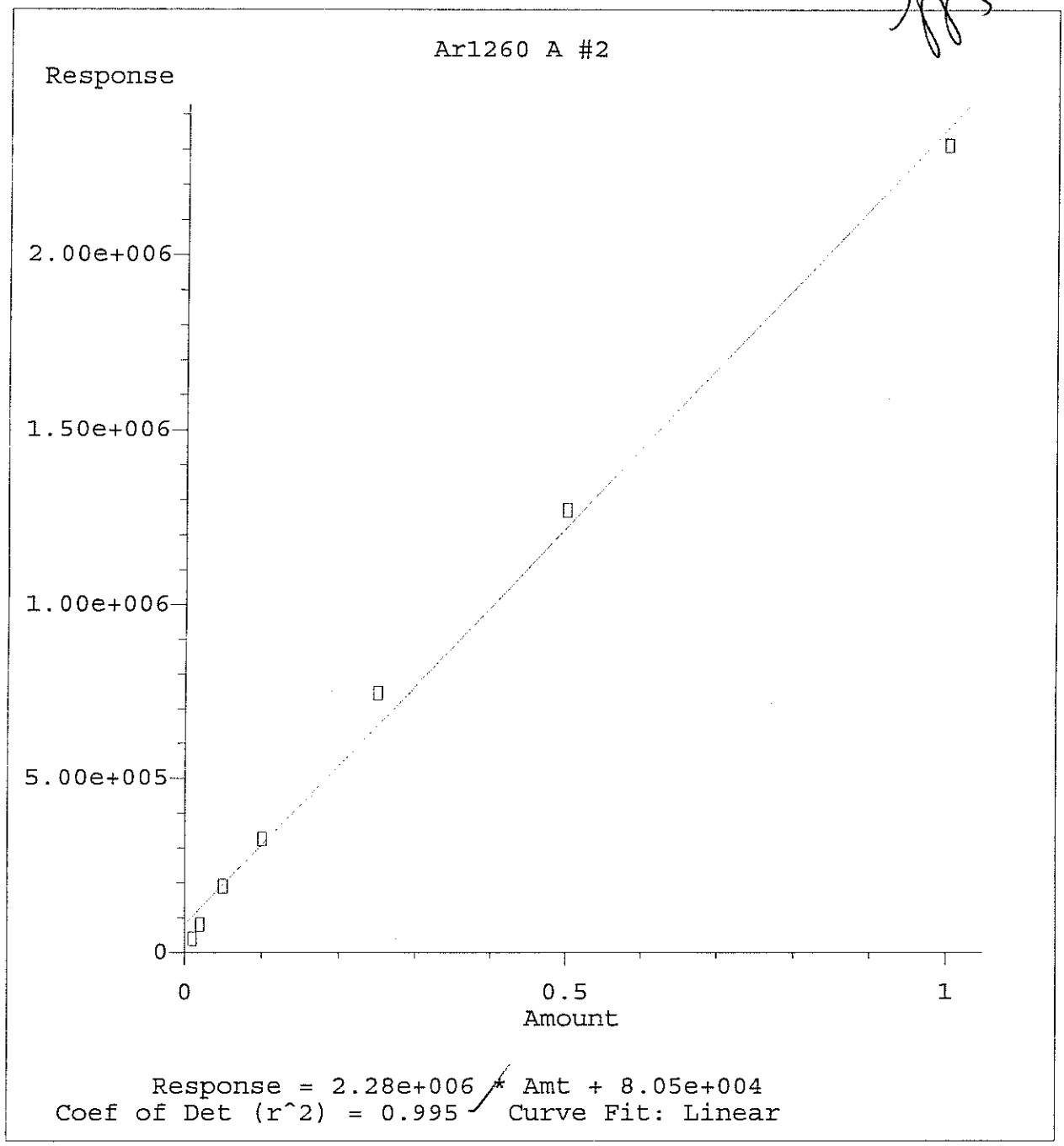
Method Name: C:\HPCHEM\2\METHODS\PBQ05166.M
Calibration Table Last Updated: Wed May 17 10:06:32 2006

5.17.06



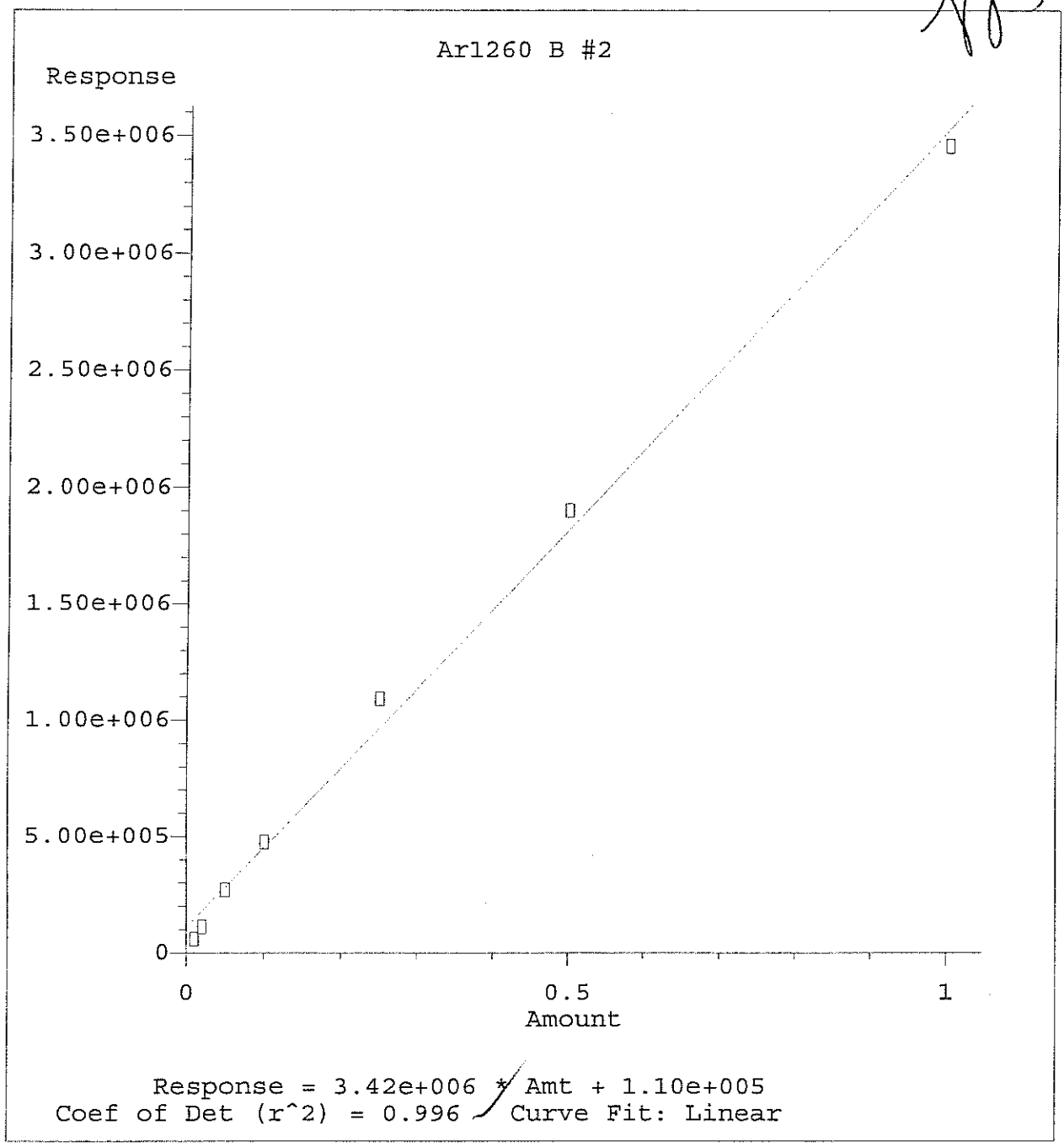
Method Name: C:\HPCHEM\2\METHODS\PBQ05166.M
Calibration Table Last Updated: Wed May 17 10:06:32 2006

5.17.06



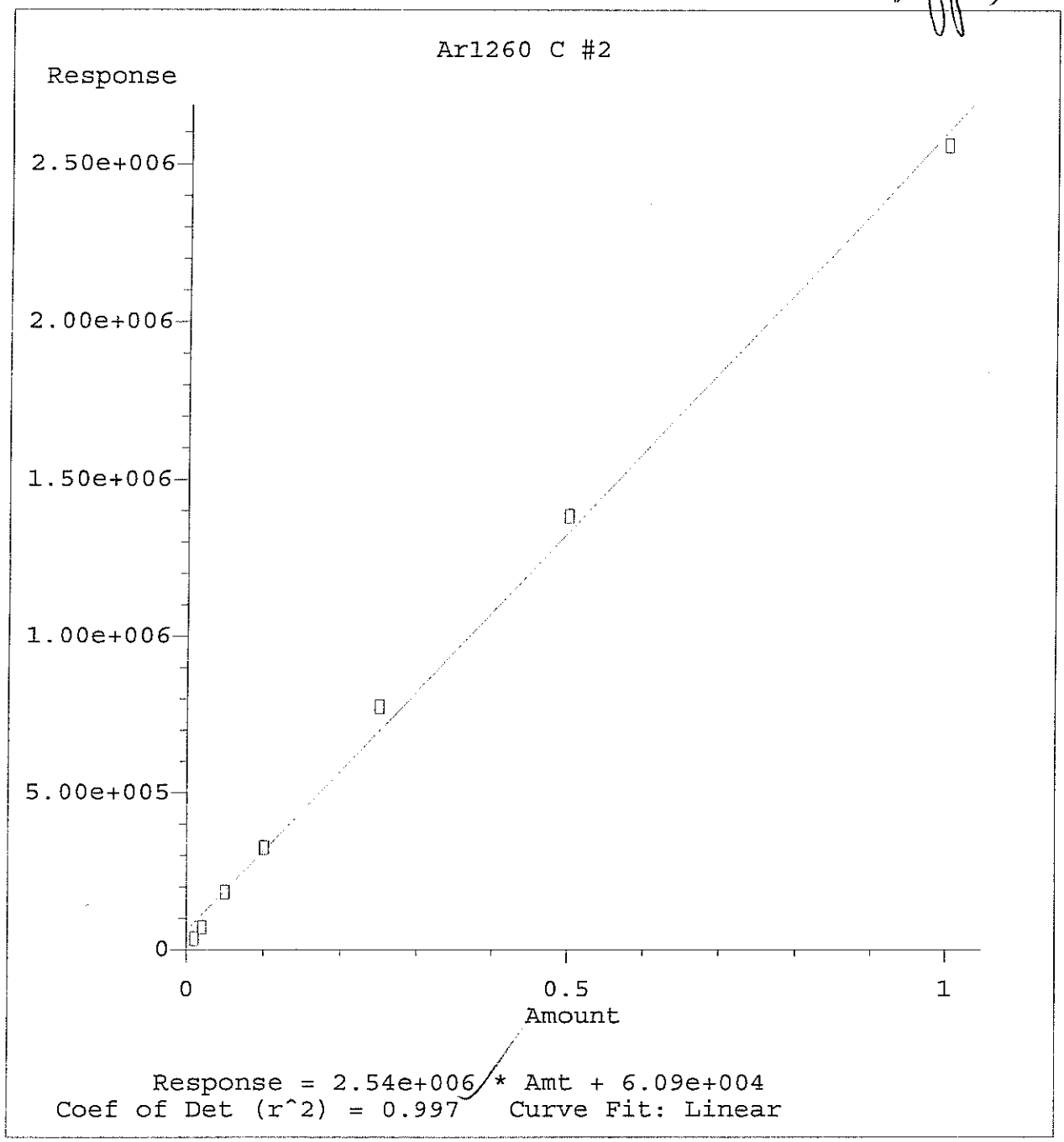
Method Name: C:\HPCHEM\2\METHODS\PBQ05166.M
Calibration Table Last Updated: Wed May 17 10:06:32 2006

5.17.06



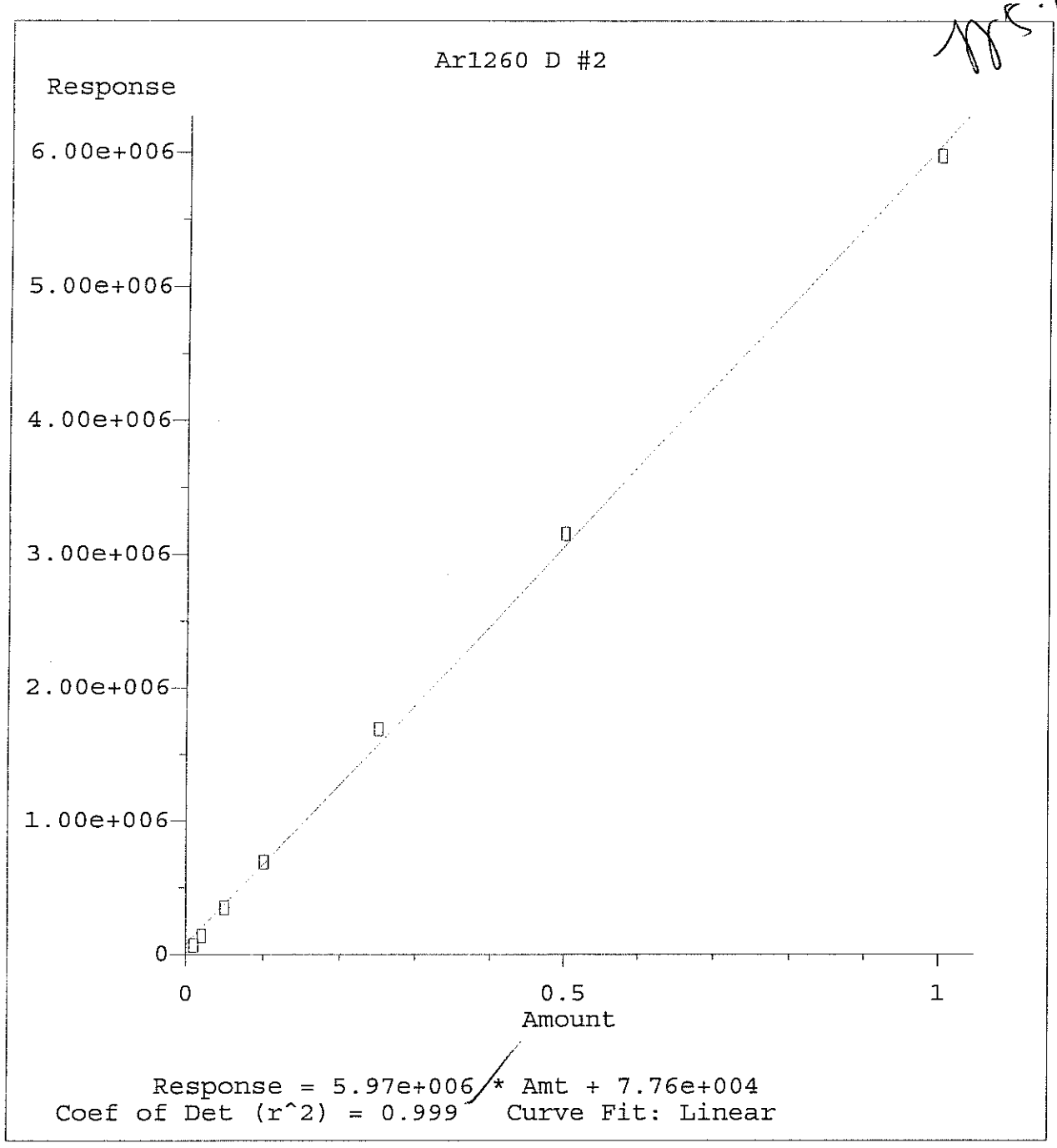
Method Name: C:\HPCHEM\2\METHODS\PBQ05166.M
Calibration Table Last Updated: Wed May 17 10:06:32 2006

5.17.06



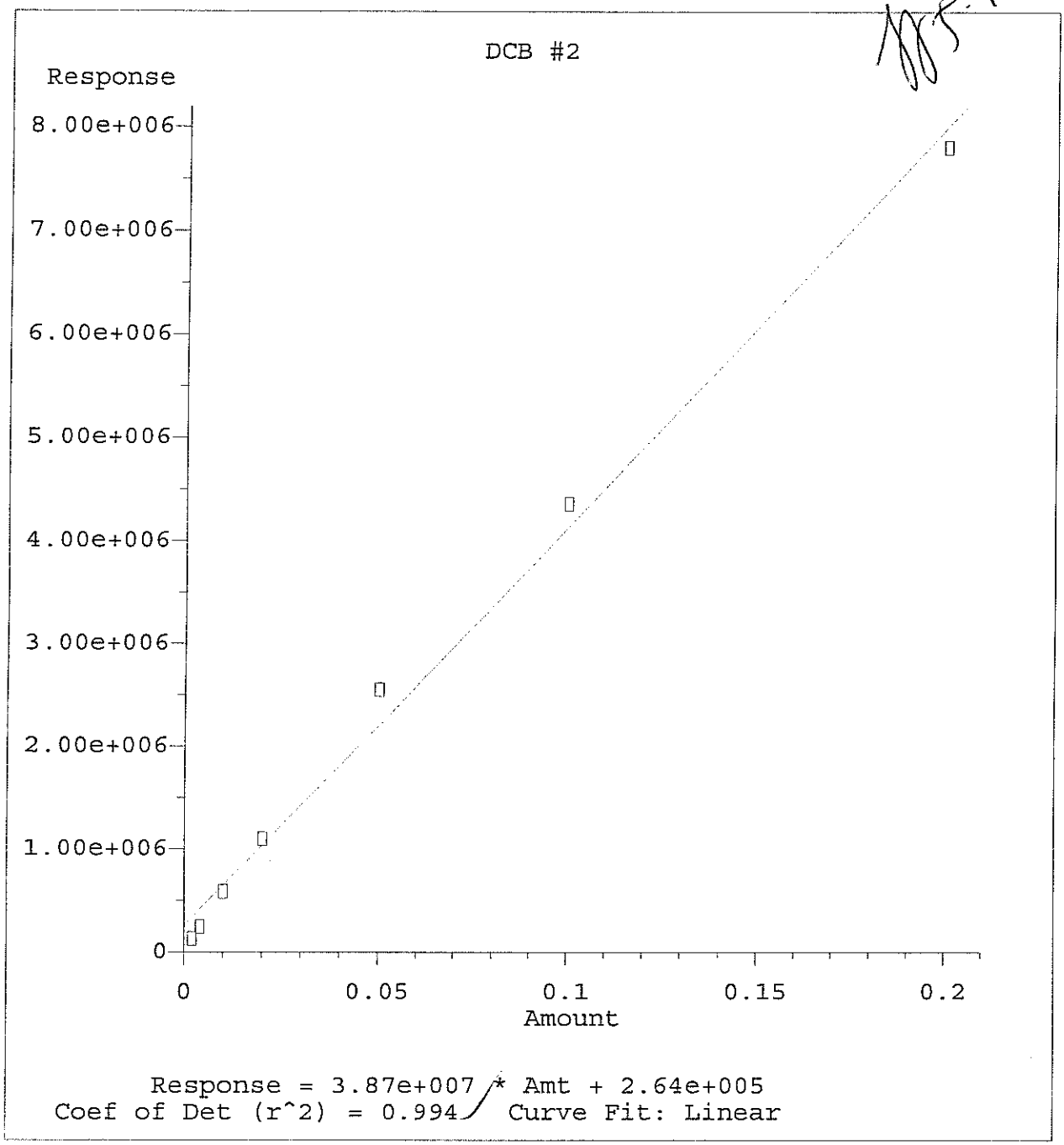
Method Name: C:\HPCHEM\2\METHODS\PBQ05166.M
Calibration Table Last Updated: Wed May 17 10:06:32 2006

MS-17-06



Method Name: C:\HPCHEM\2\METHODS\PBQ05166.M
Calibration Table Last Updated: Wed May 17 10:06:32 2006

5-17-06



Method Name: C:\HPCHEM\2\METHODS\PBQ05166.M
Calibration Table Last Updated: Wed May 17 10:06:32 2006

Quantitation Report (Not Reviewed)

Signal #1 : C:\HPCHEM\2\DATA\051606-F\F55868.D\ECD1A.CH Vial: 20
Signal #2 : C:\HPCHEM\2\DATA\051606-F\F55868.D\ECD2B.CH
Acq On : 16 May 2006 9:35 pm Operator:
Sample : P4024 16/60 0.5 CC Inst : Instrumen
Misc : Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 17 10:08 2006 Quant Results File: PBQ05166.RES

Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 10:06:32 2006
Response via : Initial Calibration
DataAcq Meth : PSTQUICK.M

5.17.06 5.17.06

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film

Compound	RT#1	RT#2	Resp#1	Resp#2	ug/ml	ug/ml
System Monitoring Compounds						
1) S TCX	2.64	2.72	3626028	3966484	0.104	0.105
Spiked Amount	0.100	Range 46 - 122		Recovery =	104.00%	105.00%
9) S DCB	10.39	10.04	3115536	4423256	0.104	0.107
Spiked Amount	0.100	Range 26 - 86		Recovery =	104.00%#	107.00%#
Target Compounds						
2) L1 Ar1016 A	3.43	3.37	832389	883397	0.488	0.496
3) L1 Ar1016 B	3.92	3.86	1587624	2003833	0.487	0.500
4) L1 Ar1016 C	4.09	4.02	656707	761021	0.483	0.501
Sum Ar1016 A			3076721	3648252	1.459	1.496
Average Ar1016 A					0.486 97%	0.499 100%
5) L2 Ar1260 A	6.66	6.35	1060994	1236509	0.507	0.507
6) L2 Ar1260 B	7.03	6.82	1233068	1866651	0.493	0.514
7) L2 Ar1260 C	7.92	7.53	954369	1330389	0.482	0.499
8) L2 Ar1260 D	8.49	8.59	934025	2982100	0.481	0.486
Sum Ar1260 A			4182456	7415649	1.963	2.006
Average Ar1260 A					0.491 99%	0.501 100%

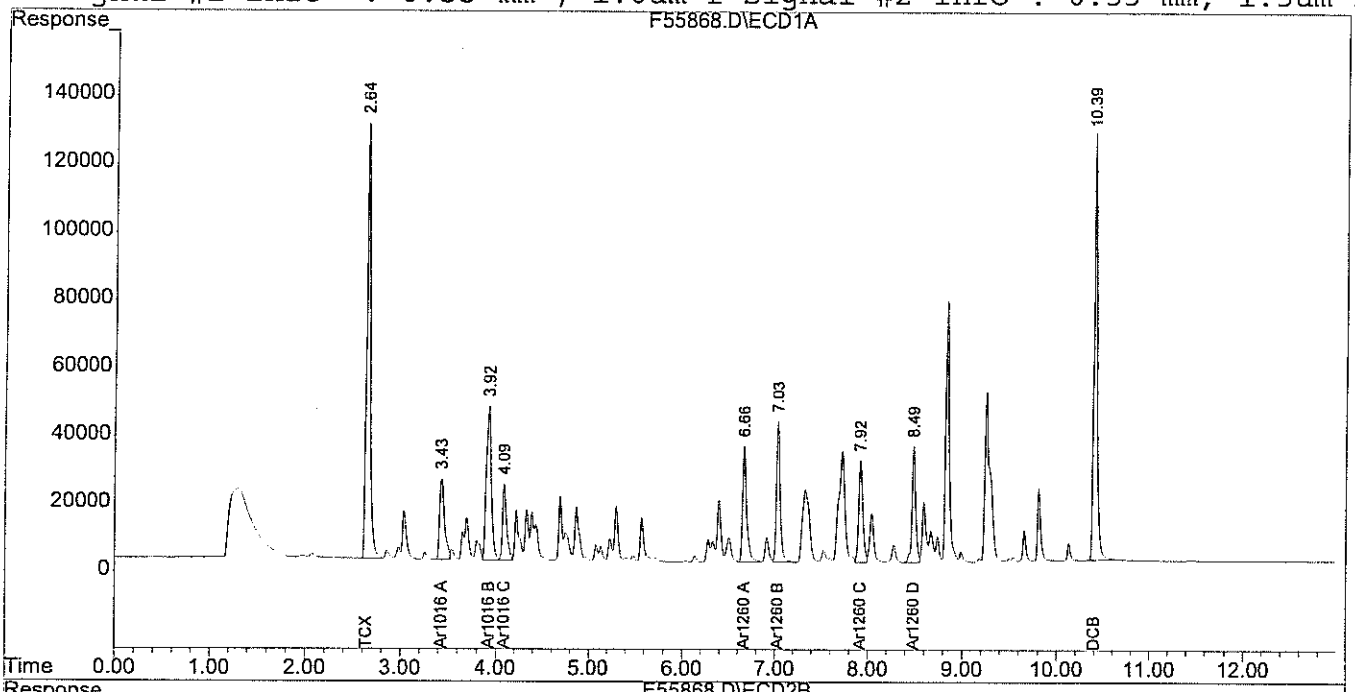
(f)=RT Delta > 1/2 Window (#)=Amounts differ by > 25% (m)=manual int

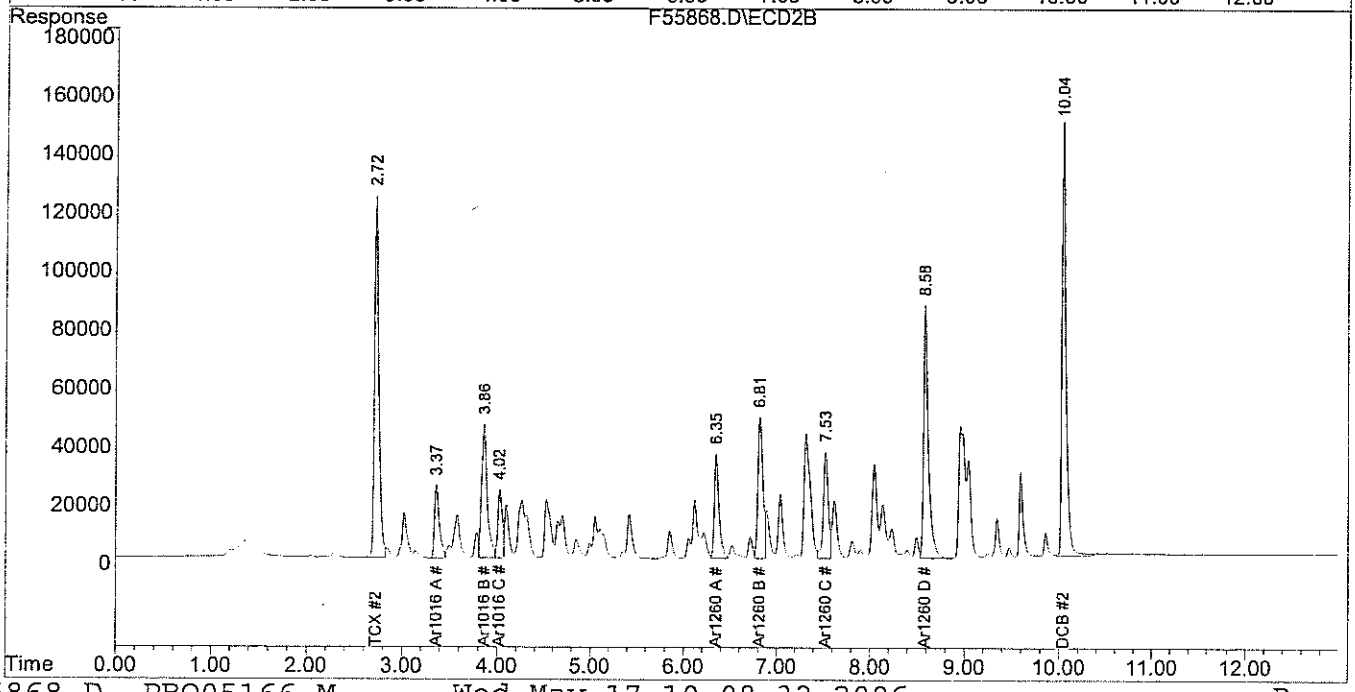
Quantitation Report (Not Reviewed)

Signal #1 : C:\HPCHEM\2\DATA\051606-F\F55868.D\ECD1A.CH Vial: 20
Signal #2 : C:\HPCHEM\2\DATA\051606-F\F55868.D\ECD2B.CH
Acq On : 16 May 2006 9:35 pm Operator:
Sample : P4024 16/60 0.5 CC Inst : Instrumen
Misc : Multiplr: 1.00
IntFile Signal #1: PCBINT.E IntFile Signal #2: PCBINT2.E
Quant Time: May 17 10:08 2006 Quant Results File: PBQ05166.RES

Quant Method : C:\HPCHEM\2\METHODS\PBQ05166.M (Chemstation Integrator)
Title : Aroclor 1016/1260
Last Update : Wed May 17 10:06:32 2006
Response via : Multiple Level Calibration
DataAcq Meth : PSTQUICK.M

Volume Inj. : 3 ul
Signal #1 Phase : DB-1701 Widebore Signal #2 Phase: DB-5 Widebore
Signal #1 Info : 0.53 mm , 1.0um f Signal #2 Info : 0.53 mm, 1.5um film





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7A
PCB CONTINUING CALIBRATION CHECK

Site ID: ME

Project No.: 218736.002 GC Column #1: RTX-CLP Pesticides I Data Package ID: 56637

Instrument ID: F GC Column #2: RTX-CLP Pesticides II ID: 0.53 mm

File #: F56151 Cal. Check Date: 05/30/2006 12:30

COMPOUND	True Value (ug/mL)	Measured Value (ug/mL)	% Recovery	*	Limits %R	Total Out
PCB 1016	0.50	0.473	95		85-115	0
PCB1260	0.50	0.460	92		85-115	0
PCB 1016#2	0.50	0.486	97		85-115	0
PCB 1260 #2	0.50	0.490	98		85-115	0

7A
PCB CONTINUING CALIBRATION CHECK

Site ID: ME

Project No.: 218736.002 GC Column #1: RTX-CLP Pesticides I Data Package ID: 56637

Instrument ID: F GC Column #2: RTX-CLP Pesticides II ID: 0.53 mm

File #: F56166 Cal. Check Date: 05/30/2006 18:36

COMPOUND	True Value (ug/mL)	Measured Value (ug/mL)	% Recovery	*	Limits %R	Total Out
PCB 1016	0.50	0.440	88		85-115	0
PCB1260	0.50	0.429	86		85-115	0
PCB 1016#2	0.50	0.467	93		85-115	0
PCB 1260 #2	0.50	0.445	89		85-115	0

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7A
PCB CONTINUING CALIBRATION CHECK

Site ID: ME

Project No.: 218736.002 GC Column #1: RTX-CLP Pesticides I Data Package ID: 56637

Instrument ID: F GC Column #2: RTX-CLP Pesticides II ID: 0.53 mm

File #: F56218 Cal. Check Date: 05/31/2006 9:43

COMPOUND	True Value (ug/mL)	Measured Value (ug/mL)	% Recovery	*	Limits %R	Total Out
PCB 1016	0.25	0.258	103		85-115	0
PCB1260	0.25	0.247	99		85-115	0
PCB 1016#2	0.25	0.260	104		85-115	0
PCB 1260 #2	0.25	0.268	107		85-115	0

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7A
PCB CONTINUING CALIBRATION CHECK

Site ID: ME

Project No.: 218736.002 GC Column #1: RTX-CLP Pesticides I Data Package ID: 56637

Instrument ID: F GC Column #2: RTX-CLP Pesticides II ID: 0.53 mm

File #: F56221 Cal. Check Date: 05/31/2006 10:27

COMPOUND	True Value (ug/mL)	Measured Value (ug/mL)	% Recovery	*	Limits %R	Total Out
PCB 1016	0.50	0.503	101		85-115	0
PCB1260	0.50	0.491	98		85-115	0
PCB 1016#2	0.50	0.505	101		85-115	0
PCB 1260 #2	0.50	0.413	83		85-115	0

PCB SYSTEM MONITORING COMPOUND RETENTION TIME SUMMARY

Quote#: SDG: 56637

Project No: 218736.002 Site ID: GP Old Town Mill Location: me

GC Columns: RTX-CLPesticidesI ID: 0.32(mm)

Instrument ID: F RTX-CLPesticidesII

Date Analyzed: 5/16/06

Lab File ID (Standard) : F55866 Time Analyzed: 20:51

					COLUMN #1		COLUMN #1	
					SMC1	SMC 2	SMC1	SMC 2
	PESTICIDE STANDARD				2.64	10.39	2.73	10.04
	UPPER LIMIT				2.69	10.44	2.78	10.09
	LOWER LIMIT				2.59	10.34	2.68	9.99
	SAMPLE ID	LAB FILE ID	DATE	TIME	COLUMN #1 SMC1 #	COLUMN #1 SMC 2 #	COLUMN #2 SMC1 #	COLUMN #2 SMC 2 #
01	P4016 16/60 0.01	F55861.D	05/16/06	19:38	2.64	10.39	2.72	10.04
02	P4017 16/60 0.02	F55862.D	05/16/06	19:53	2.64	10.40	2.73	10.04
03	P4018 16/60 0.05	F55863.D	05/16/06	20:07	2.64	10.40	2.73	10.05
04	P4019 16/60 0.1	F55864.D	05/16/06	20:22	2.64	10.40	2.73	10.05
05	P4020 16/60 0.25	F55865.D	05/16/06	20:36	2.64	10.40	2.73	10.04
06	P4021 16/60 0.5	F55866.D	05/16/06	20:51	2.64	10.39	2.73	10.04
07	P4022 16/60 1.0	F55867.D	05/16/06	21:20	2.64	10.40	2.73	10.04
08	P4024 16/60 0.5 CC	F55868.D	05/16/06	21:35	2.64	10.39	2.72	10.04
09	P3883 AR 1221	F55869.D	05/16/06	21:50	NA	NA	NA	NA
10	P3884 AR 1232	F55870.D	05/16/06	22:04	NA	NA	NA	NA
11	P3885 AR 1242	F55871.D	05/16/06	22:19	NA	NA	NA	NA
12	P3886 AR 1248	F55872.D	05/16/06	22:34	NA	NA	NA	NA
13	P3994 AR 1254	F55873.D	05/16/06	22:48	NA	NA	NA	NA
14	P3887 AR 1262	F55874.D	05/16/06	23:03	NA	NA	NA	NA
15	P3888 AR 1268	F55875.D	05/16/06	23:17	NA	NA	NA	NA
16	P4057 AR 1254 0.0	F55876.D	05/16/06	23:47	NA	NA	NA	NA
17	P4056 AR 1254 0.0	F55877.D	05/16/06	0:01	NA	NA	NA	NA
18	P4055 AR 1254 0.1	F55878.D	05/17/06	0:16	NA	NA	NA	NA
19	P4054 AR 1254 0.2	F55879.D	05/17/06	0:31	NA	NA	NA	NA
20	P4053 AR 1254 0.5	F55880.D	05/17/06	0:45	NA	NA	NA	NA
21	P4052 AR 1254 1.0	F55881.D	05/17/06	0:59	NA	NA	NA	NA
22	P4010 AR 1254 0.5	F55882.D	05/17/06	1:13	2.64	10.37	2.72	10.02
23	ZZZZZ	F55883.D	5/17/06	1:28	2.64	10.37	2.72	10.02
24	ZZZZZ	F55885.D	5/17/06	2:12	2.64	10.37	2.72	10.02
25	ZZZZZ	F55886.D	5/17/06	2:26	2.64	10.37	2.72	10.02
26	ZZZZZ	F55895.D	5/17/06	4:53	2.64	10.38	2.73	10.03
27	ZZZZZ	F55896.D	5/17/06	5:07	2.64	10.39	2.73	10.03
28	ZZZZZ	F55897.D	5/17/06	5:22	2.64	10.39	2.73	10.03
29	P4021 16/60 0.5	F55900.D	5/17/06	6:06	2.64	10.39	2.72	10.04
30	P4053 AR 1254 0.5	F55901.D	5/17/06	6:21	NA	NA	NA	NA
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PCB SYSTEM MONITORING COMPOUND RETENTION TIME SUMMARY

Quote#: SDG: 56637

Project No: 218736.002 Site ID: GP Old Town Mill Location: ME

GC Columns: RTX-CLPesticidesI ID: 0.32(mm)

Instrument ID: F RTX-CLPesticidesII

Date Analyzed: 5/30/06

Lab File ID (Standard) : F56151 Time Analyzed: 12:30

		COLUMN #1		COLUMN #1	
		SMC1	SMC 2	SMC1	SMC 2
	PESTICIDE STANDARD	2.64	10.42	2.73	10.07
	UPPER LIMIT	2.69	10.47	2.78	10.12
	LOWER LIMIT	2.59	10.37	2.68	10.02

	SAMPLE ID	LAB FILE ID	DATE	TIME	COLUMN #1 SMC1 #	COLUMN #1 SMC 2 #	COLUMN #2 SMC1 #	COLUMN #2 SMC 2 #
01	P4021 16/60 0.5	F56151.D	05/30/06	12:30	2.64	10.42	2.73	10.07
02	B05306PAS, A/C	F56152.D	05/30/06	14:27	2.70 *	10.49 *	2.77	10.13 *
03	L05306PASB, A/C	F56153.D	05/30/06	14:42	2.65	10.44	2.74	10.09
04	LD05306PASB, A/C	F56154.D	05/30/06	14:56	2.65	10.43	2.73	10.07
05	56637-4, A/C	F56155.D	05/30/06	15:11	2.64	10.42	2.73	10.07
06	56637-6, A/C	F56156.D	05/30/06	15:26	2.64	10.42	2.73	10.07
07	56637-7, A/C	F56157.D	05/30/06	16:10	2.69	10.46	2.76	10.10
08	56637-8, A/C	F56158.D	05/30/06	16:24	2.65	10.43	2.73	10.07
09	56637-9, A/C	F56159.D	05/30/06	16:39	2.64	10.41	2.73	10.05
10	56637-10, A/C	F56160.D	05/30/06	16:53	2.64	10.41	2.72	10.05
11	56637-10, MS, A/C	F56161.D	05/30/06	17:23	2.64	10.40	2.72	10.05
12	56637-10, MSD, A/C	F56162.D	05/30/06	17:37	2.64	10.40	2.72	10.05
13	56637-12, A/C	F56163.D	05/30/06	17:52	2.64	10.39	2.72	10.04
14	56637-13, A/C	F56164.D	05/30/06	18:07	2.64	10.40	2.72	10.05
15	ZZZZZ	F56165.D	05/30/06	18:21	2.64	10.41	2.72	10.05
16	P4021 16/60 0.5	F56166.D	05/30/06	18:36	2.64	10.40	2.72	10.05
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PCB SYSTEM MONITORING COMPOUND RETENTION TIME SUMMARY

Quote#: SDG: 56637

Project No: 218736.002 Site ID: GP Old Town Mill Location: ME

GC Columns: RTX-CLPesticidesI ID: 0.32(mm)

Instrument ID: F RTX-CLPesticidesII

Date Analyzed: 5/31/06

Lab File ID (Standard) : F56151 Time Analyzed: 9:43

		COLUMN #1		COLUMN #1	
		SMC1	SMC 2	SMC1	SMC 2
	PESTICIDE STANDARD	2.63	10.40	2.72	10.05
	UPPER LIMIT	2.68	10.45	2.77	10.10
	LOWER LIMIT	2.58	10.35	2.67	10.00

	SAMPLE ID	LAB FILE ID	DATE	TIME	COLUMN #1 SMC1 #	COLUMN #1 SMC 2 #	COLUMN #2 SMC1 #	COLUMN #2 SMC 2 #
01	P4020 16/60 0.25	F56218.D	05/31/06	9:43	2.63	10.40	2.72	10.05
02	B05306PAS, RR A	F56219.D	05/31/06	9:58	2.67	10.43	2.75	10.07
03	56637-19, 1:20	F56220.D	05/31/06	14:42	2.64	10.41	2.72	10.05
04	P4021 16/60 0.5	F56221.D	05/31/06	10:27	2.64	10.42	2.72	10.06
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Maine Environmental Laboratory

Report of Analyses

One Main Street Yarmouth, Maine 04096-1107 Tel (207) 846-6569 Fax (207) 846-9066 e-mail: melab@ime.net

Steve Knollmeyer
Analytics Environmental Lab, LLC
195 Commerce Way, Suite E
Portsmouth, NH 03801

Report Date: June 01, 2006

Report No.: AEL4078-06

Enclosed are the results of the analyses requested on your samples as received by the laboratory. Samples were received in acceptable condition and analyzed within method holding times with all quality control data within laboratory acceptance limits unless noted. Reporting detection limits are the minimum levels for reporting quantitative data. These limits are 3.18 times the method detection limit as defined in CFR 40 Part 136, Appendix B. Data reported between the reporting and method detection limits are J flagged as estimated. Maine Environmental Laboratory is certified by Maine, Massachusetts, New Hampshire and NELAP (cert.#2031). A list of certified parameters is available on request. The results reported herein conform to the most current NELAP standards, where applicable, unless otherwise narrated in the body of the report. This report shall not be reproduced, except in full, without the written consent of the laboratory.

The complete report consists of the following sections: Maine Environmental Laboratory report
Chain of custody forms

Project Name/ID: GP Old Town Mill 218736.002

Date Received: 05/26/06

Sampler Name: No Data

References

EPA600/4-79-020, Methods for Chemical Analysis of Water and Wastes, USEPA, Cincinnati, Ohio, March 1983.

SW846, Test Methods for Evaluating Solid Waste, Physical/Chemical Methods, USEPA, third edition, 1986.

CLP - USEPA CLP Statement of Work for Inorganics, ILMO3.0.

Standard Methods for the Examination of Water and Wastewater, 18th edition, APHA,AWWA,WPCF, 1992.

EPA/600/R-93/100 Methods for the Determination of Inorganic Substances in Environmental Samples, Aug. 1993.

HACH - Chemical Oxygen Demand, Method 8000, Hach Handbook of Water Analysis, Hach Chemical Company, 1979.

Authorized signature

Herbert S. Kodis, laboratory director

PROJECTSITE	SAMPLE_POINT_ID	LAB_SAMPLE_ID	ANALYSIS_LAB	SAMPLE_DATE	SAMPLE_TYPE	PARAMETER_NAME	CONCENTRATION	LAB_QUALIFIER3	QUANTITATION_LIMIT	PARAMETER_UNITS	TEST	ANALYSIS_DATE	ANALYSIS_TIME	PREP_METHOD	WEIGHT_B	MDL	SAMPLE_FILTER	Ti
10 Text	30 Text	30 Text	4 Text	Date	2 Text	100 Text	Number (Double	3 Text	Number (Double)	8 Text	15 Text	Date	Time	15 Text	3 Text	Number (Double	2 Text	
1	2	3	4	5	6	8	9	10	11	12	13	14	15	18	20	23	34	3
GP OLD TOWN MILL	SB-5(0-4')(56637-10)	AEL407806-01	ME	05/25/06	SL	ARSENIC	7.0		0.4	MG/KG	SW7060A	05/30/06		SW3050B	DRY	0.1		
GP OLD TOWN MILL	SB-5(0-4')(56637-10)	AEL407806-01	ME	05/25/06	SL	BARIUM	20	J	40	MG/KG	SW7080A	05/31/06		SW3050B	DRY	10		
GP OLD TOWN MILL	SB-5(0-4')(56637-10)	AEL407806-01	ME	05/25/06	SL	CADMIUM		U	2.1	MG/KG	SW7130	05/31/06		SW3050B	DRY	0.5		
GP OLD TOWN MILL	SB-5(0-4')(56637-10)	AEL407806-01	ME	05/25/06	SL	CHROMIUM	19		8	MG/KG	SW7190	05/31/06		SW3050B	DRY	2		
GP OLD TOWN MILL	SB-5(0-4')(56637-10)	AEL407806-01	ME	05/25/06	SL	LEAD	16		11	MG/KG	SW7420	05/31/06		SW3050B	DRY	3		
GP OLD TOWN MILL	SB-5(0-4')(56637-10)	AEL407806-01	ME	05/25/06	SL	MERCURY		U	0.05	MG/KG	SW7471A	05/30/06			DRY	0.02		
GP OLD TOWN MILL	SB-5(0-4')(56637-10)	AEL407806-01	ME	05/25/06	SL	SELENIUM	0.5		0.4	MG/KG	SW7740	05/30/06		SW3050B	DRY	0.1		
GP OLD TOWN MILL	SB-5(0-4')(56637-10)	AEL407806-01	ME	05/25/06	SL	SILVER		U	8	MG/KG	SW7760A	05/31/06		SW3050B	DRY	2		
GP OLD TOWN MILL	SB-6(0-4')(56637-11)	AEL407806-02	ME	05/25/06	SL	ARSENIC	6.1		0.4	MG/KG	SW7060A	05/30/06		SW3050B	DRY	0.1		
GP OLD TOWN MILL	SB-6(0-4')(56637-11)	AEL407806-02	ME	05/25/06	SL	BARIUM	40		40	MG/KG	SW7080A	05/31/06		SW3050B	DRY	10		
GP OLD TOWN MILL	SB-6(0-4')(56637-11)	AEL407806-02	ME	05/25/06	SL	CADMIUM		U	2.0	MG/KG	SW7130	05/31/06		SW3050B	DRY	0.5		
GP OLD TOWN MILL	SB-6(0-4')(56637-11)	AEL407806-02	ME	05/25/06	SL	CHROMIUM	22		7	MG/KG	SW7190	05/31/06		SW3050B	DRY	2		
GP OLD TOWN MILL	SB-6(0-4')(56637-11)	AEL407806-02	ME	05/25/06	SL	LEAD	12		10	MG/KG	SW7420	05/31/06		SW3050B	DRY	3		
GP OLD TOWN MILL	SB-6(0-4')(56637-11)	AEL407806-02	ME	05/25/06	SL	MERCURY		U	0.07	MG/KG	SW7471A	05/30/06			DRY	0.02		
GP OLD TOWN MILL	SB-6(0-4')(56637-11)	AEL407806-02	ME	05/25/06	SL	SELENIUM	0.1	J	0.4	MG/KG	SW7740	05/30/06		SW3050B	DRY	0.1		
GP OLD TOWN MILL	SB-6(0-4')(56637-11)	AEL407806-02	ME	05/25/06	SL	SILVER		U	7	MG/KG	SW7760A	05/31/06		SW3050B	DRY	2		
GP OLD TOWN MILL	SB-15(0-4')(56637-18)	AEL407806-03	ME	05/25/06	SL	ARSENIC	3.2		0.5	MG/KG	SW7060A	05/30/06		SW3050B	DRY	0.1		
GP OLD TOWN MILL	SB-15(0-4')(56637-18)	AEL407806-03	ME	05/25/06	SL	BARIUM	50		50	MG/KG	SW7080A	05/31/06		SW3050B	DRY	10		
GP OLD TOWN MILL	SB-15(0-4')(56637-18)	AEL407806-03	ME	05/25/06	SL	CADMIUM		U	2.3	MG/KG	SW7130	05/31/06		SW3050B	DRY	0.6		
GP OLD TOWN MILL	SB-15(0-4')(56637-18)	AEL407806-03	ME	05/25/06	SL	CHROMIUM	17		8	MG/KG	SW7190	05/31/06		SW3050B	DRY	2		
GP OLD TOWN MILL	SB-15(0-4')(56637-18)	AEL407806-03	ME	05/25/06	SL	LEAD	6	J	11	MG/KG	SW7420	05/31/06		SW3050B	DRY	3		
GP OLD TOWN MILL	SB-15(0-4')(56637-18)	AEL407806-03	ME	05/25/06	SL	MERCURY		U	0.06	MG/KG	SW7471A	05/30/06			DRY	0.02		
GP OLD TOWN MILL	SB-15(0-4')(56637-18)	AEL407806-03	ME	05/25/06	SL	SELENIUM	0.2	J	0.5	MG/KG	SW7740	05/30/06		SW3050B	DRY	0.1		
GP OLD TOWN MILL	SB-15(0-4')(56637-18)	AEL407806-03	ME	05/25/06	SL	SILVER		U	8	MG/KG	SW7760A	05/31/06		SW3050B	DRY	2		

463/486

PROJECT/SITE	SAMPLE_POINT_ID	LAB_SAMPLE_ID	ANALYSIS_LAB	SAMPLE_DATE	SAMPLE_TYPE	PARAMETER_NAME	CONCENTRATION	LAB_QUALIFIERS	QUANTITATION_LIMIT	PARAMETER_UNITS	TEST	ANALYSIS_DATE	ANALYSIS_TIME	PREP_METHOD	WEIGHT_B	MDL	SAMPLE_FILTER	SAMPLE_TYPE
10 Text	30 Text	30 Text	4 Text	Date	2 Text	100 Text	Number (Double	3 Text	Number (Double)	8 Text	15 Text	Date	Time	15 Text	3 Text	Number (Double	2 Text	Ti
1	2	3	4	5	6	8	9	10	11	12	13	14	15	18	20	23	34	3
GP OLD TOWN MILL	SB-21(0-4')(56637-9)	AEL407806-04	ME	05/25/06	SL	ARSENIC	6.4		0.4	MG/KG	SW7060A	05/30/06		SW3050B	DRY	0.1		
GP OLD TOWN MILL	SB-21(0-4')(56637-9)	AEL407806-04	ME	05/25/06	SL	BARIUM	30	J	40	MG/KG	SW7080A	05/31/06		SW3050B	DRY	10		
GP OLD TOWN MILL	SB-21(0-4')(56637-9)	AEL407806-04	ME	05/25/06	SL	CADMIUM		U	2.1	MG/KG	SW7130	05/31/06		SW3050B	DRY	0.5		
GP OLD TOWN MILL	SB-21(0-4')(56637-9)	AEL407806-04	ME	05/25/06	SL	CHROMIUM	23		7	MG/KG	SW7190	05/31/06		SW3050B	DRY	2		
GP OLD TOWN MILL	SB-21(0-4')(56637-9)	AEL407806-04	ME	05/25/06	SL	LEAD	5	J	10	MG/KG	SW7420	05/31/06		SW3050B	DRY	3		
GP OLD TOWN MILL	SB-21(0-4')(56637-9)	AEL407806-04	ME	05/25/06	SL	MERCURY		U	0.05	MG/KG	SW7471A	05/30/06			DRY	0.02		
GP OLD TOWN MILL	SB-21(0-4')(56637-9)	AEL407806-04	ME	05/25/06	SL	SELENIUM	0.1	J	0.4	MG/KG	SW7740	05/30/06		SW3050B	DRY	0.1		
GP OLD TOWN MILL	SB-21(0-4')(56637-9)	AEL407806-04	ME	05/25/06	SL	SILVER		U	7	MG/KG	SW7760A	05/31/06		SW3050B	DRY	2		
GP OLD TOWN MILL	SB-22(0-4')(56637-4)	AEL407806-05	ME	05/25/06	SL	ARSENIC	7.5		0.5	MG/KG	SW7060A	05/30/06		SW3050B	DRY	0.1		
GP OLD TOWN MILL	SB-22(0-4')(56637-4)	AEL407806-05	ME	05/25/06	SL	BARIUM	20	J	50	MG/KG	SW7080A	05/31/06		SW3050B	DRY	10		
GP OLD TOWN MILL	SB-22(0-4')(56637-4)	AEL407806-05	ME	05/25/06	SL	CADMIUM	0.7	J	2.4	MG/KG	SW7130	05/31/06		SW3050B	DRY	0.6		
GP OLD TOWN MILL	SB-22(0-4')(56637-4)	AEL407806-05	ME	05/25/06	SL	CHROMIUM	106		8	MG/KG	SW7190	05/31/06		SW3050B	DRY	2		
GP OLD TOWN MILL	SB-22(0-4')(56637-4)	AEL407806-05	ME	05/25/06	SL	LEAD	171		12	MG/KG	SW7420	05/31/06		SW3050B	DRY	4		
GP OLD TOWN MILL	SB-22(0-4')(56637-4)	AEL407806-05	ME	05/25/06	SL	MERCURY	0.02	J	0.06	MG/KG	SW7471A	05/30/06			DRY	0.02		
GP OLD TOWN MILL	SB-22(0-4')(56637-4)	AEL407806-05	ME	05/25/06	SL	SELENIUM	0.7		0.5	MG/KG	SW7740	05/30/06		SW3050B	DRY	0.1		
GP OLD TOWN MILL	SB-22(0-4')(56637-4)	AEL407806-05	ME	05/25/06	SL	SILVER		U	8	MG/KG	SW7760A	05/31/06		SW3050B	DRY	2		
GP OLD TOWN MILL	SB-23(0-4')(56637-12)	AEL407806-06	ME	05/25/06	SL	ARSENIC	5.3		0.4	MG/KG	SW7060A	05/30/06		SW3050B	DRY	0.1		
GP OLD TOWN MILL	SB-23(0-4')(56637-12)	AEL407806-06	ME	05/25/06	SL	BARIUM	40		40	MG/KG	SW7080A	05/31/06		SW3050B	DRY	10		
GP OLD TOWN MILL	SB-23(0-4')(56637-12)	AEL407806-06	ME	05/25/06	SL	CADMIUM		U	2.2	MG/KG	SW7130	05/31/06		SW3050B	DRY	0.6		
GP OLD TOWN MILL	SB-23(0-4')(56637-12)	AEL407806-06	ME	05/25/06	SL	CHROMIUM	28		8	MG/KG	SW7190	05/31/06		SW3050B	DRY	2		
GP OLD TOWN MILL	SB-23(0-4')(56637-12)	AEL407806-06	ME	05/25/06	SL	LEAD	7	J	11	MG/KG	SW7420	05/31/06		SW3050B	DRY	3		
GP OLD TOWN MILL	SB-23(0-4')(56637-12)	AEL407806-06	ME	05/25/06	SL	MERCURY		U	0.06	MG/KG	SW7471A	05/30/06			DRY	0.02		
GP OLD TOWN MILL	SB-23(0-4')(56637-12)	AEL407806-06	ME	05/25/06	SL	SELENIUM		U	0.4	MG/KG	SW7740	05/30/06		SW3050B	DRY	0.1		
GP OLD TOWN MILL	SB-23(0-4')(56637-12)	AEL407806-06	ME	05/25/06	SL	SILVER		U	8	MG/KG	SW7760A	05/31/06		SW3050B	DRY	2		

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PROJECT/SITE	SAMPLE_POINT_ID	LAB_SAMPLE_ID	ANALYSIS_LAB	SAMPLE_DATE	SAMPLE_TYPE	PARAMETER_NAME	CONCENTRATION	LAB_QUALIFIERS	QUANTITATION_LIMIT	PARAMETER_UNITS	TEST	ANALYSIS_DATE	ANALYSIS_TIME	PREP_METHOD	WEIGHT_B	MDL	SAMPLE_FILTER	Ti
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1	2	3	4	5	6	8	9	10	11	12	13	14	15	18	20	23	34	3
GP OLD TOWN MILL	SB-30(0-4')(56637-15)	AEL407806-07	ME	05/25/06	SL	ARSENIC	5.7		0.5	MG/KG	SW7060A	05/30/06		SW3050B	DRY	0.1		
GP OLD TOWN MILL	SB-30(0-4')(56637-15)	AEL407806-07	ME	05/25/06	SL	BARIUM	160		50	MG/KG	SW7080A	05/31/06		SW3050B	DRY	10		
GP OLD TOWN MILL	SB-30(0-4')(56637-15)	AEL407806-07	ME	05/25/06	SL	CADMIUM		U	2.5	MG/KG	SW7130	05/31/06		SW3050B	DRY	0.6		
GP OLD TOWN MILL	SB-30(0-4')(56637-15)	AEL407806-07	ME	05/25/06	SL	CHROMIUM	12		9	MG/KG	SW7190	05/31/06		SW3050B	DRY	3		
GP OLD TOWN MILL	SB-30(0-4')(56637-15)	AEL407806-07	ME	05/25/06	SL	LEAD	129		12	MG/KG	SW7420	05/31/06		SW3050B	DRY	4		
GP OLD TOWN MILL	SB-30(0-4')(56637-15)	AEL407806-07	ME	05/25/06	SL	MERCURY	0.09		0.06	MG/KG	SW7471A	05/30/06			DRY	0.02		
GP OLD TOWN MILL	SB-30(0-4')(56637-15)	AEL407806-07	ME	05/25/06	SL	SELENIUM	0.2	J	0.5	MG/KG	SW7740	05/30/06		SW3050B	DRY	0.1		
GP OLD TOWN MILL	SB-30(0-4')(56637-15)	AEL407806-07	ME	05/25/06	SL	SILVER		U	9	MG/KG	SW7760A	05/31/06		SW3050B	DRY	3		
GP OLD TOWN MILL	SB-31(0-4')(56637-14)	AEL407806-08	ME	05/25/06	SL	ARSENIC	9.0		0.6	MG/KG	SW7060A	05/30/06		SW3050B	DRY	0.2		
GP OLD TOWN MILL	SB-31(0-4')(56637-14)	AEL407806-08	ME	05/25/06	SL	BARIUM	130		60	MG/KG	SW7080A	05/31/06		SW3050B	DRY	20		
GP OLD TOWN MILL	SB-31(0-4')(56637-14)	AEL407806-08	ME	05/25/06	SL	CADMIUM		U	3.2	MG/KG	SW7130	05/31/06		SW3050B	DRY	0.8		
GP OLD TOWN MILL	SB-31(0-4')(56637-14)	AEL407806-08	ME	05/25/06	SL	CHROMIUM	28		11	MG/KG	SW7190	05/31/06		SW3050B	DRY	3		
GP OLD TOWN MILL	SB-31(0-4')(56637-14)	AEL407806-08	ME	05/25/06	SL	LEAD	96		16	MG/KG	SW7420	05/31/06		SW3050B	DRY	5		
GP OLD TOWN MILL	SB-31(0-4')(56637-14)	AEL407806-08	ME	05/25/06	SL	MERCURY	0.52		0.08	MG/KG	SW7471A	05/30/06			DRY	0.03		
GP OLD TOWN MILL	SB-31(0-4')(56637-14)	AEL407806-08	ME	05/25/06	SL	SELENIUM	0.2	J	0.6	MG/KG	SW7740	05/30/06		SW3050B	DRY	0.2		
GP OLD TOWN MILL	SB-31(0-4')(56637-14)	AEL407806-08	ME	05/25/06	SL	SILVER		U	11	MG/KG	SW7760A	05/31/06		SW3050B	DRY	3		
GP OLD TOWN MILL	SB-37(0-4')(56637-17)	AEL407806-09	ME	05/25/06	SL	ARSENIC	5.2		0.4	MG/KG	SW7060A	05/30/06		SW3050B	DRY	0.1		
GP OLD TOWN MILL	SB-37(0-4')(56637-17)	AEL407806-09	ME	05/25/06	SL	BARIUM	20	J	40	MG/KG	SW7080A	05/31/06		SW3050B	DRY	10		
GP OLD TOWN MILL	SB-37(0-4')(56637-17)	AEL407806-09	ME	05/25/06	SL	CADMIUM		U	2.2	MG/KG	SW7130	05/31/06		SW3050B	DRY	0.5		
GP OLD TOWN MILL	SB-37(0-4')(56637-17)	AEL407806-09	ME	05/25/06	SL	CHROMIUM	18		8	MG/KG	SW7190	05/31/06		SW3050B	DRY	2		
GP OLD TOWN MILL	SB-37(0-4')(56637-17)	AEL407806-09	ME	05/25/06	SL	LEAD	10	J	11	MG/KG	SW7420	05/31/06		SW3050B	DRY	3		
GP OLD TOWN MILL	SB-37(0-4')(56637-17)	AEL407806-09	ME	05/25/06	SL	MERCURY		U	0.05	MG/KG	SW7471A	05/30/06			DRY	0.02		
GP OLD TOWN MILL	SB-37(0-4')(56637-17)	AEL407806-09	ME	05/25/06	SL	SELENIUM		U	0.4	MG/KG	SW7740	05/30/06		SW3050B	DRY	0.1		
GP OLD TOWN MILL	SB-37(0-4')(56637-17)	AEL407806-09	ME	05/25/06	SL	SILVER		U	8	MG/KG	SW7760A	05/31/06		SW3050B	DRY	2		

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1	2	3	4	5	6	8	9	10	11	12	13	14	15	18	20	23	34	3
GP OLD TOWN MILL	SB-38(0-4')(56637-16)	AEL407806-10	ME	05/25/06	SL	ARSENIC	6.9		0.4	MG/KG	SW7060A	05/30/06		SW3050B	DRY	0.1		
GP OLD TOWN MILL	SB-38(0-4')(56637-16)	AEL407806-10	ME	05/25/06	SL	BARIUM	40		40	MG/KG	SW7080A	05/31/06		SW3050B	DRY	10		
GP OLD TOWN MILL	SB-38(0-4')(56637-16)	AEL407806-10	ME	05/25/06	SL	CADMIUM		U	2.1	MG/KG	SW7130	05/31/06		SW3050B	DRY	0.5		
GP OLD TOWN MILL	SB-38(0-4')(56637-16)	AEL407806-10	ME	05/25/06	SL	CHROMIUM	26		8	MG/KG	SW7190	05/31/06		SW3050B	DRY	2		
GP OLD TOWN MILL	SB-38(0-4')(56637-16)	AEL407806-10	ME	05/25/06	SL	LEAD	18		11	MG/KG	SW7420	05/31/06		SW3050B	DRY	3		
GP OLD TOWN MILL	SB-38(0-4')(56637-16)	AEL407806-10	ME	05/25/06	SL	MERCURY	0.04	J	0.05	MG/KG	SW7471A	05/30/06			DRY	0.02		
GP OLD TOWN MILL	SB-38(0-4')(56637-16)	AEL407806-10	ME	05/25/06	SL	SELENIUM		U	0.4	MG/KG	SW7740	05/30/06		SW3050B	DRY	0.1		
GP OLD TOWN MILL	SB-38(0-4')(56637-16)	AEL407806-10	ME	05/25/06	SL	SILVER		U	8	MG/KG	SW7760A	05/31/06		SW3050B	DRY	2		
GP OLD TOWN MILL	SB-41(0-4')(56637-19)	AEL407806-11	ME	05/25/06	SL	ARSENIC	5.1		0.4	MG/KG	SW7060A	05/30/06		SW3050B	DRY	0.1		
GP OLD TOWN MILL	SB-41(0-4')(56637-19)	AEL407806-11	ME	05/25/06	SL	BARIUM	10	J	40	MG/KG	SW7080A	05/31/06		SW3050B	DRY	10		
GP OLD TOWN MILL	SB-41(0-4')(56637-19)	AEL407806-11	ME	05/25/06	SL	CADMIUM		U	2.2	MG/KG	SW7130	05/31/06		SW3050B	DRY	0.6		
GP OLD TOWN MILL	SB-41(0-4')(56637-19)	AEL407806-11	ME	05/25/06	SL	CHROMIUM	14		8	MG/KG	SW7190	05/31/06		SW3050B	DRY	2		
GP OLD TOWN MILL	SB-41(0-4')(56637-19)	AEL407806-11	ME	05/25/06	SL	LEAD	15		11	MG/KG	SW7420	05/31/06		SW3050B	DRY	3		
GP OLD TOWN MILL	SB-41(0-4')(56637-19)	AEL407806-11	ME	05/25/06	SL	MERCURY	0.04	J	0.05	MG/KG	SW7471A	05/30/06			DRY	0.02		
GP OLD TOWN MILL	SB-41(0-4')(56637-19)	AEL407806-11	ME	05/25/06	SL	SELENIUM		U	0.4	MG/KG	SW7740	05/30/06		SW3050B	DRY	0.1		
GP OLD TOWN MILL	SB-41(0-4')(56637-19)	AEL407806-11	ME	05/25/06	SL	SILVER		U	8	MG/KG	SW7760A	05/31/06		SW3050B	DRY	2		
GP OLD TOWN MILL	PILE 1 (56637-22)	AEL407806-12	ME	05/26/06	SL	ARSENIC	5.8		0.4	MG/KG	SW7060A	05/30/06		SW3050B	DRY	0.1		
GP OLD TOWN MILL	PILE 1 (56637-22)	AEL407806-12	ME	05/26/06	SL	BARIUM	10	J	40	MG/KG	SW7080A	05/31/06		SW3050B	DRY	10		
GP OLD TOWN MILL	PILE 1 (56637-22)	AEL407806-12	ME	05/26/06	SL	CADMIUM		U	2.2	MG/KG	SW7130	05/31/06		SW3050B	DRY	0.6		
GP OLD TOWN MILL	PILE 1 (56637-22)	AEL407806-12	ME	05/26/06	SL	CHROMIUM	21		8	MG/KG	SW7190	05/31/06		SW3050B	DRY	2		
GP OLD TOWN MILL	PILE 1 (56637-22)	AEL407806-12	ME	05/26/06	SL	LEAD	12		11	MG/KG	SW7420	05/31/06		SW3050B	DRY	3		
GP OLD TOWN MILL	PILE 1 (56637-22)	AEL407806-12	ME	05/26/06	SL	MERCURY		U	0.06	MG/KG	SW7471A	05/30/06			DRY	0.02		
GP OLD TOWN MILL	PILE 1 (56637-22)	AEL407806-12	ME	05/26/06	SL	SELENIUM		U	0.4	MG/KG	SW7740	05/30/06		SW3050B	DRY	0.1		
GP OLD TOWN MILL	PILE 1 (56637-22)	AEL407806-12	ME	05/26/06	SL	SILVER		U	8	MG/KG	SW7760A	05/31/06		SW3050B	DRY	2		

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10 Text	30 Text	30 Text	4 Text	Date	2 Text	100 Text	Number (Double	3 Text	Number (Double)	8 Text	15 Text	Date	Time	15 Text	3 Text	Number (Double	2 Text	
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GP OLD TOWN MILL	PILE 2 (56637-23)	AEL407806-13	ME	05/26/06	SL	ARSENIC	6.0		0.4	MG/KG	SW7060A	05/30/06		SW3050B	DRY	0.1		
GP OLD TOWN MILL	PILE 2 (56637-23)	AEL407806-13	ME	05/26/06	SL	BARIUM		U	40	MG/KG	SW7080A	05/31/06		SW3050B	DRY	10		
GP OLD TOWN MILL	PILE 2 (56637-23)	AEL407806-13	ME	05/26/06	SL	CADMIUM		U	2.1	MG/KG	SW7130	05/31/06		SW3050B	DRY	0.5		
GP OLD TOWN MILL	PILE 2 (56637-23)	AEL407806-13	ME	05/26/06	SL	CHROMIUM	22		7	MG/KG	SW7190	05/31/06		SW3050B	DRY	2		
GP OLD TOWN MILL	PILE 2 (56637-23)	AEL407806-13	ME	05/26/06	SL	LEAD	8	J	10	MG/KG	SW7420	05/31/06		SW3050B	DRY	3		
GP OLD TOWN MILL	PILE 2 (56637-23)	AEL407806-13	ME	05/26/06	SL	MERCURY		U	0.05	MG/KG	SW7471A	05/30/06			DRY	0.02		
GP OLD TOWN MILL	PILE 2 (56637-23)	AEL407806-13	ME	05/26/06	SL	SELENIUM		U	0.4	MG/KG	SW7740	05/30/06		SW3050B	DRY	0.1		
GP OLD TOWN MILL	PILE 2 (56637-23)	AEL407806-13	ME	05/26/06	SL	SILVER		U	7	MG/KG	SW7760A	05/31/06		SW3050B	DRY	2		
GP OLD TOWN MILL	SB-15 (56637-20)	AEL407806-14	ME	05/26/06	GW	ARSENIC	0.016		0.004	MG/L	SW7060A	05/30/06				0.001	FM	
GP OLD TOWN MILL	SB-15 (56637-20)	AEL407806-14	ME	05/26/06	GW	BARIUM		U	0.4	MG/L	SW7080A	05/30/06		SW3005A		0.1	FM	
GP OLD TOWN MILL	SB-15 (56637-20)	AEL407806-14	ME	05/26/06	GW	CADMIUM		U	0.0006	MG/L	SW7131A	05/27/06		SW3020A		0.0002	FM	
GP OLD TOWN MILL	SB-15 (56637-20)	AEL407806-14	ME	05/26/06	GW	CHROMIUM		U	0.005	MG/L	SW7191	05/31/06		SW3020A		0.002	FM	
GP OLD TOWN MILL	SB-15 (56637-20)	AEL407806-14	ME	05/26/06	GW	LEAD		U	0.006	MG/L	SW7421	05/31/06		SW3020A		0.002	FM	
GP OLD TOWN MILL	SB-15 (56637-20)	AEL407806-14	ME	05/26/06	GW	MERCURY		U	0.0005	MG/L	SW7470A	05/30/06				0.0002	FM	
GP OLD TOWN MILL	SB-15 (56637-20)	AEL407806-14	ME	05/26/06	GW	SELENIUM	0.004	J	0.006	MG/L	SW7740	05/30/06				0.002	FM	
GP OLD TOWN MILL	SB-15 (56637-20)	AEL407806-14	ME	05/26/06	GW	SILVER		U	0.05	MG/L	SW7760A	05/30/06				0.02	FM	
GP OLD TOWN MILL	SB-37 (56637-3)	AEL407806-15	ME	05/26/06	GW	ARSENIC	0.002	J	0.004	MG/L	SW7060A	05/30/06				0.001	FM	
GP OLD TOWN MILL	SB-37 (56637-3)	AEL407806-15	ME	05/26/06	GW	BARIUM		U	0.4	MG/L	SW7080A	05/30/06		SW3005A		0.1	FM	
GP OLD TOWN MILL	SB-37 (56637-3)	AEL407806-15	ME	05/26/06	GW	CADMIUM		U	0.0006	MG/L	SW7131A	05/27/06		SW3020A		0.0002	FM	
GP OLD TOWN MILL	SB-37 (56637-3)	AEL407806-15	ME	05/26/06	GW	CHROMIUM		U	0.005	MG/L	SW7191	05/31/06		SW3020A		0.002	FM	
GP OLD TOWN MILL	SB-37 (56637-3)	AEL407806-15	ME	05/26/06	GW	LEAD		U	0.006	MG/L	SW7421	05/31/06		SW3020A		0.002	FM	
GP OLD TOWN MILL	SB-37 (56637-3)	AEL407806-15	ME	05/26/06	GW	MERCURY	0.0007		0.0005	MG/L	SW7470A	05/30/06				0.0002	FM	
GP OLD TOWN MILL	SB-37 (56637-3)	AEL407806-15	ME	05/26/06	GW	SELENIUM	0.002	J	0.006	MG/L	SW7740	05/30/06				0.002	FM	
GP OLD TOWN MILL	SB-37 (56637-3)	AEL407806-15	ME	05/26/06	GW	SILVER		U	0.05	MG/L	SW7760A	05/30/06				0.02	FM	

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1	2	3	4	5	6	8	9	10	11	12	13	14	15	18	20	23	34	3
GP OLD TOWN MILL	MW-5 (56637-1)	AEL407806-16	ME	05/26/06	GW	ARSENIC		U	0.004	MG/L	SW7060A	05/30/06				0.001	FM	
GP OLD TOWN MILL	MW-5 (56637-1)	AEL407806-16	ME	05/26/06	GW	BARIUM		U	0.4	MG/L	SW7080A	05/30/06		SW3005A		0.1	FM	
GP OLD TOWN MILL	MW-5 (56637-1)	AEL407806-16	ME	05/26/06	GW	CADMIUM		U	0.0006	MG/L	SW7131A	05/27/06		SW3020A		0.0002	FM	
GP OLD TOWN MILL	MW-5 (56637-1)	AEL407806-16	ME	05/26/06	GW	CHROMIUM		U	0.005	MG/L	SW7191	05/31/06		SW3020A		0.002	FM	
GP OLD TOWN MILL	MW-5 (56637-1)	AEL407806-16	ME	05/26/06	GW	LEAD		U	0.006	MG/L	SW7421	05/31/06		SW3020A		0.002	FM	
GP OLD TOWN MILL	MW-5 (56637-1)	AEL407806-16	ME	05/26/06	GW	MERCURY	0.0002	J	0.0005	MG/L	SW7470A	05/30/06				0.0002	FM	
GP OLD TOWN MILL	MW-5 (56637-1)	AEL407806-16	ME	05/26/06	GW	SELENIUM	0.004	J	0.006	MG/L	SW7740	05/30/06				0.002	FM	
GP OLD TOWN MILL	MW-5 (56637-1)	AEL407806-16	ME	05/26/06	GW	SILVER		U	0.05	MG/L	SW7760A	05/30/06				0.02	FM	
GP OLD TOWN MILL	MW-3 (56637-21)	AEL407806-17	ME	05/26/06	GW	ARSENIC	0.003	J	0.004	MG/L	SW7060A	05/30/06				0.001	FM	
GP OLD TOWN MILL	MW-3 (56637-21)	AEL407806-17	ME	05/26/06	GW	BARIUM		U	0.4	MG/L	SW7080A	05/30/06		SW3005A		0.1	FM	
GP OLD TOWN MILL	MW-3 (56637-21)	AEL407806-17	ME	05/26/06	GW	CADMIUM		U	0.0006	MG/L	SW7131A	05/27/06		SW3020A		0.0002	FM	
GP OLD TOWN MILL	MW-3 (56637-21)	AEL407806-17	ME	05/26/06	GW	CHROMIUM		U	0.005	MG/L	SW7191	05/31/06		SW3020A		0.002	FM	
GP OLD TOWN MILL	MW-3 (56637-21)	AEL407806-17	ME	05/26/06	GW	LEAD		U	0.006	MG/L	SW7421	05/31/06		SW3020A		0.002	FM	
GP OLD TOWN MILL	MW-3 (56637-21)	AEL407806-17	ME	05/26/06	GW	MERCURY	0.0002	J	0.0005	MG/L	SW7470A	05/30/06				0.0002	FM	
GP OLD TOWN MILL	MW-3 (56637-21)	AEL407806-17	ME	05/26/06	GW	SELENIUM		U	0.006	MG/L	SW7740	05/30/06				0.002	FM	
GP OLD TOWN MILL	MW-3 (56637-21)	AEL407806-17	ME	05/26/06	GW	SILVER		U	0.05	MG/L	SW7760A	05/30/06				0.02	FM	
GP OLD TOWN MILL	MW-4 (56637-2)	AEL407806-18	ME	05/26/06	GW	ARSENIC	0.003	J	0.004	MG/L	SW7060A	05/30/06				0.001	FM	
GP OLD TOWN MILL	MW-4 (56637-2)	AEL407806-18	ME	05/26/06	GW	BARIUM		U	0.4	MG/L	SW7080A	05/30/06		SW3005A		0.1	FM	
GP OLD TOWN MILL	MW-4 (56637-2)	AEL407806-18	ME	05/26/06	GW	CADMIUM		U	0.0006	MG/L	SW7131A	05/27/06		SW3020A		0.0002	FM	
GP OLD TOWN MILL	MW-4 (56637-2)	AEL407806-18	ME	05/26/06	GW	CHROMIUM		U	0.005	MG/L	SW7191	05/31/06		SW3020A		0.002	FM	
GP OLD TOWN MILL	MW-4 (56637-2)	AEL407806-18	ME	05/26/06	GW	LEAD		U	0.006	MG/L	SW7421	05/31/06		SW3020A		0.002	FM	
GP OLD TOWN MILL	MW-4 (56637-2)	AEL407806-18	ME	05/26/06	GW	MERCURY		U	0.0005	MG/L	SW7470A	05/30/06				0.0002	FM	
GP OLD TOWN MILL	MW-4 (56637-2)	AEL407806-18	ME	05/26/06	GW	SELENIUM	0.004	J	0.006	MG/L	SW7740	05/30/06				0.002	FM	
GP OLD TOWN MILL	MW-4 (56637-2)	AEL407806-18	ME	05/26/06	GW	SILVER		U	0.05	MG/L	SW7760A	05/30/06				0.02	FM	

Katahdin ANALYTICAL SERVICES
340 County Road No. 5
P.O. Box 720
Westbrook, ME 04092
Tel: (207) 874-2400
Fax: (207) 775-4029

Analytics

CHAIN of CUSTODY

PLEASE BEAR DOWN AND PRINT LEGIBLY IN PEN

Page ____ of ____

Client: Woodard + Curran
Contact: Karl Kasper
Phone #: (207) 774-2112
Fax #: (207) 774-6635

Address: 41 Hutchins Drive
City: Portland
State: ME
Zip Code: 04102

Purchase Order #: 218736.002
Proj. Name / No.:
Katahdin Quote #:

Bill (if different than above):
Address:

Sampler (Print / Sign): Josh Bowe
Copies To:

LAB USE ONLY
WORK ORDER #:
KATAHDIN PROJECT NUMBER ____

REMARKS: ____

SHIPPING INFO: ☐ FED EX ☐ UPS ☐ CLIENT
AIRBILL NO: ____
TEMP°C ____ ☐ TEMP BLANK ☐ INTACT ☐ NOT INTACT

ANALYSIS AND CONTAINER TYPE PRESERVATIVES

*	Sample Description	Date / Time coll'd	Matrix	No. of Cntrs.	Filt. ☐Y ☑N 40ml 8260 B Unpreserved	Filt. ☐Y ☑N PCB Unpreserved 4oz	Filt. ☐Y ☑N Metals RCRA8 Unpreserved 8oz	Filt. ☐Y ☑N 40ml low soil H2O preserved	Filt. ☐Y ☐N 40ml med soil MeOH preserved	Filt. ☐Y ☐N SVOA (8270) Unpreserved 2oz	Filt. ☐Y ☐N SVOA (8270) Unpreserved 4oz	Filt. ☐Y ☐N	Filt. ☐Y ☐N	Filt. ☐Y ☐N
	SB-1 (6-10')	5/25 / 8:00	GW	2	2									
	SB-21 (4-8')	/ 11:50	Soil	1		1								
	SB-35 (0-6")	/ 7:00	Soil	1		1								
	SB-5 (4-6')	/ 13:20	Soil	1		1								
4	SB-21 (0-4')	/ 11:40	Soil	6		1	1	2	1	1				
1	SB-5 (0-4')	/ 13:15	Soil	6		1	1	2	1	1				
2	SB-6 (0-4')	/ 14:00	Soil	5			1	2	1		1			
6	SB-23 (0-4)'	/ 15:30	Soil	3		1	1				1			
	SB-24 (0-4)'	/ 14:10	Soil	1		1								
		/												
		/												
		/												
		/												
		/												
		/												
		/												

4 Soil

COMMENTS

Relinquished By: (Signature)	Date / Time	Received By: (Signature)	Relinquished By: (Signature)	Date / Time	Received By: (Signature)
Josh Bowe	5/25 16:16	[signature]			

Maine Environmental Laboratory - Quality Control Data Report

Report: ATL 4078-06 **Sample matrix:** SOLID **Date:** 6/1/16

Duplicate Samples - sample specific quality control

Parameter	Sample ID	data-mg/kg	duplicate	% RSD	Acceptance limits
HG	4078-01 Rep	ND	0.02	N/A	Duplicate acceptance limits are ± 35% at concentrations 5 times the RDL.
HG	4078-12	ND	ND	N/A	
Se	4078-01	0.5	0.4	N/A	
Se	4078-12	ND	ND	N/A	
As	4078-01	7.0	5.5	24.0%	
As	4078-12	5.8	6.2	6.7%	
Pb	4078-10	16	18	11.8%	
Cr	4078-10	26	25	3.9%	
Ag	4078-10	ND	ND	N/A	
Ba	4078-10	40	40	0.0%	
Cd	4078-10	ND	ND	N/A	

Matrix Spike Samples - sample specific quality control - accept. limits: **75% - 125%** @ 4X spike

Parameter	Sample ID	data-mg/kg (wet wet)	spike addition	spike data	% Recovery
HG	4078-05	0.02	0.50	0.54	104%
HG	4078-10	0.04	0.50	0.58	108%
Se	4078-05	0.6	1.0	1.8	120%
Se	4078-10	0.0	1.0	1.0	100%
As	4078-05	6.3	(w) +/- 5.0	11.0	94%
As	4078-10	6.5	5.0	10.8	86%
Pb	4078-01	15	50	64	98%
Cr	4078-01	18	20	39	105%
Ag	4078-01	ND(0)	25	24	96%
Ba	4078-01	20	200	240	110%
Cd	4078-01	0.0 ND	5.0	4.6	92%

Maine Environmental Laboratory - Quality Control Data Report

Report: ALL 4078-06 **Sample matrix:** SOLID **Date:** 6/1/16

Laboratory Control Samples - analytical batch quality control

Parameter	LCS - ID	data-mg/kg	True Value	% Recovery	Acceptance limits
Hg	0.50 Hg LCS	0.55	0.50	110%	Laboratory control samples acceptance limits are **80% - 120%** recovery.
Se	5-0 Se LCS	4.6	5-0	92%	
As	4.0 Se LCS	3.9	4-0	98%	
Pb	ASIT-3	67	67	100%	
Cr	ASIT-3	11	11	100%	
Ag	WM-2	17	18	94%	
Ba	WM-1	410	450	91%	
Cd	WM-2	3.9	4.0	98%	

Method Blanks - analytical batch quality control

Parameter	report data	data-mg/kg	RDL	Acceptance limits
Hg	ND	0.00	0.05	Data for method blanks must be **less than or equal to the RDL** (reporting quantitation detection limit).
Se	ND	0.0	0.4	
As	ND	0.0	0.4	
Pb	ND	0	10	
Cr	ND	0	7	
Ag	ND	0	7	
Ba	ND	0	40	
Cd	ND	0.0	2.0	

Maine Environmental Laboratory - Quality Control Data Report

Report: AL 4078-06 **Sample matrix:** AQ **Date:** 01/16

Duplicate Samples - sample specific quality control

Parameter	Sample ID	data -mg/L	duplicate	% RSD	Acceptance limits
Hg	4078-18	0.0000	0.0000	N/A	Duplicate acceptance limits are ± 20% at concentrations 5 times the RDL.
Se	4078-16	0.004	0.005	N/A	
As	4078-16	0.000	0.000	N/A	
Ba	4078-16	0.0	0.0	N/A	
Ag	4078-16	0.00	0.00	N/A	
Cd	4078-16	0.001	0.0000	N/A	
Pb	4078-16	0.001	0.000	N/A	
Cr	4078-16	0.01	0.001	N/A	

Matrix Spike Samples - sample specific quality control - accept. limits: **75% - 125%** @ 4X spike

Parameter	Sample ID	data -mg/L	spike addition	spike data	% Recovery
Hg	4078-14	0.0000	0.0020	0.0020	100%
Se	4078-14	0.004	0.025	0.023	76%
As	4078-14	0.016	0.025	0.039	92%
Ba	4078-14	0.0	2.0	2.1	105%
Ag	4078-14	0.00	0.20	0.21	105%
Cd	4078-14	0.0000	0.0020	0.0021	105%
Pb	4078-14	0.001	0.025	0.023	88%
Cr	4078-14	0.000	0.025	0.022	88%

Maine Environmental Laboratory - Quality Control Data Report

Report: AEL 4678-06 **Sample matrix:** AG **Date:** 6/1/16

Laboratory Control Samples - analytical batch quality control

Parameter	LCS - ID	data -mg/L	True Value	% Recovery	Acceptance limits
Hg	5.0 Hg LCS	0.0054	0.0050	108%	Laboratory control samples acceptance limits are **90% - 110%** recovery.
Se	100 Se LCS	0.099	0.100	99%	
As	50 As LCS	0.051	0.050	102%	
Ba	2.0 Ba LCS	2.1	2.0	105%	
Ag	1.00 Ag LCS	0.99	1.00	99%	
Cd	3.0 Cd LCS	0.031	0.030	103%	
Pb	20 Pb LCS	0.019	0.020	95%	
Cr	25 Cr LCS	0.024	0.025	96%	

Method Blanks - analytical batch quality control

Parameter	report data	data -mg/L	RDL	Acceptance limits
Hg	ND	0.0000	0.0005	Data for method blanks must be less than or equal to the RDL (reporting quantitation detection limit).
Se	ND	0.000	0.006	
As	ND	0.000	0.004	
Ba	ND	0.0	0.4	
Ag	ND	0.00	0.05	
Cd	ND	0.0000	0.0006	
Pb	ND	0.000	0.006	
Cr	ND	0.000	0.005	

Chain Of Custody Form

analytics environmental laboratory LLC

195 Commerce Way Suite E
Portsmouth, NH 03801
Phone (603) 436-5111
Fax (603) 430-2151

For Analytics Use Only Rev. 1, 10/1/02

Samples were:
1) Shipped or hand-delivered
2) Temp blank °C 8.1°C
3) Received in good condition Y or N
4) pH checked by: _____
5) Labels checked by: _____

roject#: 218736.002 Proj. Name: GP Old Town Mill
ompany: Woodard &Curran
ontact: Karl Kasper
ddress: 41 Hutchins Drive
Portland, ME 04102
hone: (207)774-2112 PO# Quote #
ampler (Signature): [signature]

Matrix Key:
WW=Wastewater
SW=Surfacewater
GW=Groundwater
DW=Drinkingwater
S=Soil/Sludge
O=Oil
E=Extract
X=Other

Container Key
P=plastic G=glass

Preservation

Station Identification	Sample Date	Sample Time	Analysis	Degrees	4°C	HNO3	H2SO4	HCL	Methanol	Other	Matrix	Container number/type		pH	Analytics Sample #
MW-5	5/26/06	13:00	RCRA 8 Metals			✓					GW	1	500ml		
MW-5	5/26/06	13:00	SVOC 8270		✓						GW	1	1L		
MW-5	5/26/06	13:00	VOC					✓			GW	4	40ml		
MW-1	5/26/06	07:30	RCRA 8 Metals			✓					GW	1	500ml		
MW-1	5/26/06	07:30	SVOC 8270		✓						GW	1	1L		
MW-1	5/26/06	07:30	VOC					✓			GW	4	40ml		
~~MW-3~~	~~5/26/06~~	~~07:30~~													
~~MW-3~~	~~5/26/06~~	~~07:30~~													
MW-3															
~~MW-3~~			~~Sampled 5/26/06~~												
MW-4			samples 5/26/06 (Filtered ??) Dissolved?												

Received By: [signature] Time: 15:00 Date: 5/26/06
Received By: [signature] Time: 1815 Date: 5/26/06
Received By: [signature] Time: Date:

AX RESULTS? YES NO
ax # 207-774-6635
Turnaround Request
andard ☐ Priority ☐
e Date Due Date

Comments / Instructions:
Metal Sample Filtered
MISSING
~~of~~ no MW-1
~~bottles MW-3 in label~~

GW-2
XTRA MW-4
~~MW-3~~

Page 1 of 2

Relinquished By Sampler: [signature]
Relinquished By:
Relinquished By:

alytics\AEL Documents\AEL COC

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Phone (603) 436-5111
Fax (603) 430-2151

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Samples were:
1) Shipped or hand-delivered
2) Temp blank °C _____
3) Received in good condition Y or N
4) pH checked by: _____
5) Labels checked by: _____

roject#: Proj. Name:

ompany: Woodard &Curran

ontact:

ddress: 41 Hutchins Drive

Portland, ME 04102

hone: (207)774-2112 PO# Quote #

ampler (Signature):

Matrix Key:
WW=Wastewater
SW=Surfacewater
GW=Groundwater
DW=Drinkingwater
S=Soil/Sludge
O=Oil
E=Extract
X=Other

Container Key
P=plastic G=glass

Preservation: Unpres | 4°C | HNO3 | H2SO4 | HCL | Methanol | Other

Station Identification	Sample Date	Sample Time	Analysis	Matrix	Container number/type		pH	Analytics Sample #
SB-15 (0-4)	5/26/06	1310	SVOCs	GW	1	G		
SB-15 (0-4)		1330	VOCs	GW	4			
SB-15 (0-4)		1330	metals	GW	1			
MW-3		1225	SVOCs	GW	1			
MW-3		1225	VOCs	GW	3			
MW-3		1250	metals	GW	1			
Pile 1		1400	SVOCs	S	1			
Pile 1		1400	metals	S	1			
Pile 2		1410	SVOCs		1			
Pile 2		1410	Metals		1			
SB-40 (1½-5½)		0810	VOCs		2			

Comments / Instructions:

metal GW samples filtered
2-GW
2-Soil

FAX RESULTS? YES NO
fax # 207-774-6635

Turnaround Request
Standard ☐ Priority ☐
Due Date Due Date

Page___ of___

Received By: Paul Pelkus Time: 1814 Date: 5/26/06
Relinquished By Sampler: Time: Date:
Received By: Time: Date:
Relinquished By: Time: Date:
Received By: Time: Date:
Relinquished By: Time: Date:

Chain Of Custody Form

analytics environmental laboratory LLC
195 Commerce Way Suite E
Portsmouth, NH 03801
Phone (603) 436-5111
Fax (603) 430-2151

For Analytics Use Only Rev. 1, 10/1/02

Samples were:
1) Shipped or hand-delivered
2) Temp blank °C 8.1°C
3) Received in good condition Y or N
4) pH checked by: _____
5) Labels checked by: _____

Project#: 218736.002 Proj. Name: GP Old Town Mill

Company: Woodard &Curran

Contact: Karl Kasper

Address: 41 Hutchins Drive

Portland, ME 04102

Phone: (207)774-2112 PO# Quote #

Sampler (Signature):

Matrix Key:
WW=Wastewater
SW=Surfacewater
GW=Groundwater
DW=Drinkingwater
S=Soil/Sludge
O=Oil
E=Extract
X=Other

Container Key
P=plastic G=glass

Preservation

Station Identification	Sample Date	Sample Time	Analysis	Unpres	4°C	HNO3	H2SO4	HCL	Methanol	Other	Matrix	Container number/type		pH	Analytics Sample #
SB-37 GW	5/26/06	14:25	RCRA 8 Metals GW			✓					GW	1	500ml P		
SB-37		14:25	SVOC 8270									1	1L G		
SB-37		14:25	VOC					✓				4	40ml G		
✓ SB-22 (0-4)	5/25/06	1600	PCB	✓							S	1	8oz		
✓ SB-22 (0-4)	5/25/06	1600	metals	✓							S	1	4oz		
✓ SB-22 (0-4)	5/25/06	1600	SVOCs	✓							S	1	4oz		
✓ SB-22 (0-4)	5/25/06	1600	VOCs												
~~SB-39 (0-4)~~	~~5/25/06~~	~~1500~~													

Comments / Instructions: Metals filtered. (GW)

1 GW
1 - SOIL

FAX RESULTS? YES NO
Fax # 207-774-6635

Turnaround Request
Standard ☐ Priority ☒
Due Date Due Date

Received By: Time: 15:00 Date: 5/26/06
Received By: Time: 1812 Date: 5/26/06
Received By: Time: Date:

Relinquished By Sampler:
Relinquished By:
Relinquished By:

Chain Of Custody Form

analytics environmental laboratory LLC
195 Commerce Way Suite E
Portsmouth, NH 03801
Phone (603) 436-5111
Fax (603) 430-2151

For Analytics Use Only Rev. 1, 10/1/02

Samples were:
1) Shipped or hand-delivered
2) Temp blank °C _____
3) Received in good condition Y or N
4) pH checked by: _____
5) Labels checked by: _____

Project#: 218736 Proj. Name:
Company: Woodard &Curran
Contact: KARL KASPER
Address: 41 Hutchins Drive
Portland, ME 04102
Phone: (207)774-2112 PO# Quote #
Sampler (Signature): Danielle Taylor

Matrix Key:
WW=Wastewater
SW=Surfacewater
GW=Groundwater
DW=Drinkingwater
S=Soil/Sludge
O=Oil
E=Extract
X=Other

Container Key
P=plastic G=glass

Preservation: Unpres | 4°C | HNO3 | H2SO4 | HCL | Methanol | Other

Station Identification	Sample Date	Sample Time	Analysis	Matrix	Container number/type	pH	Analytics Sample #
SB-37 (0-4)	5/25/06	1340	RCRA 8 metals	S	1P 1G		
SB-15 (0-4)	5/25/06	0920	metals	S	1		
SB-15 (0-4)	5/25/06	0930	SVOCs	S	1		
SB-15 (0-4)	5/25/06	0930	VOCs	S	1		
~~SB-40 (0-4)~~	~~5/26/06~~	~~0800~~	~~metals~~	~~S~~	~~1~~		
		~~0900~~	~~SVOCs~~	~~S~~	~~1~~		
		~~0820~~	~~VOCs~~	~~S~~	~~1~~		
SB-41 (0-4)		0850	metals	S	1		
			SVOCs	S	1		
			VOCs	S	1		
			PCB	S	1		

Comments / Instructions:
3-SOIL

FAX RESULTS? YES NO
Fax # 207-774-6635

Turnaround Request
Standard ☐ Priority ☐
Due Date Due Date

Page 2 of ____

Received By:	Received By:	Received By:
Time: 16:30	Time: 1813	Time:
Date: 5/25/06	Date: 5/26/06	Date:
Relinquished By Sampler: Danielle Taylor	Relinquished By:	Relinquished By:

Chain Of Custody Form

analytics environmental laboratory LLC
195 Commerce Way Suite E
Portsmouth, NH 03801
Phone (603) 436-5111
Fax (603) 430-2151

For Analytics Use Only Rev. 1, 10/1/02

Samples were:
1) Shipped or hand-delivered
2) Temp blank °C 8.0°C
3) Received in good condition Y or N
4) pH checked by: _____
5) Labels checked by: _____

Project#: 218736 Proj. Name:
Company: Woodard &Curran
Contact: KARI KASPER
Address: 41 Hutchins Drive
Portland, ME 04102
Phone: (207)774-2112 PO# Quote #
Sampler (Signature): Danielle Taylor

Matrix Key:
WW=Wastewater
SW=Surfacewater
GW=Groundwater
DW=Drinkingwater
S=Soil/Sludge
O=Oil
E=Extract
X=Other

Container Key
P=plastic G=glass

Preservation

Station Identification	Sample Date	Sample Time	Analysis	Unpres	4°C	HNO3	H2SO4	HCL	Methanol	Other	Matrix	Container number/type		pH	Analytics Sample #
✓ SB-31 (0-4) ✓	5/25/06	10:30	8260 VOC							✓	S	2	G		
✓ "	"	"	8270 SVOC								S	1			
✓ "	"	"	RCRA 8 Metals ✓								S	1			
✓ SB-30 (0-4) ✓	"	11:25	8260 VOC									2			
✓ "	"	11:25	8270 SVOC									1			
✓ "	"	11:25	RCRA 8 Metals ✓									1			
" DT	"		DT												
✓ SB-38 (0-4)	"	13:00	8260 VOC									2			
✓ "	"	13:00	8270 SVOC									1			
✓ "	"	13:00	RCRA 8 Metals ✓									1			
✓ SB-37 (0-4)	"	13:40	8260 VOCs									2			
✓ "	"	13:40	8270 SVOC									1			

FAX RESULTS? YES NO
Fax # 207-774-6635

Turnaround Request
Standard ☐ Due Date _____
Priority ☐ Due Date _____

Comments / Instructions:
3 SOIL

Page 1 of ____

Received By:	Time: 1811	Date: 5/25/06
Received By:	Time:	Date:
Received By:	Time:	Date:
Relinquished By Sampler: Danielle Taylor		
Relinquished By:		
Relinquished By:		

Analytics\AEL Documents\AEL COC

Chain Of Custody Form

analytics environmental laboratory LLC
195 Commerce Way Suite E
Portsmouth, NH 03801
Phone (603) 436-5111
Fax (603) 430-2151

For Analytics Use Only Rev. 1, 10/1/02

Samples were:
1) Shipped or hand-delivered
2) Temp blank °C 8.1°C/4°C
3) Received in good condition Y (O) N
4) pH checked by: MEL
5) Labels checked by: MMG 5.30.06

Project#: 21873G.002 Proj. Name: GP Old Town Mill

Company: Woodard &Curran

Contact: Karl Kasper

Address: 41 Hutchins Drive

Portland, ME 04102

Phone: (207)774-2112 PO# Quote #

Sampler (Signature):

Matrix Key:
WW=Wastewater
SW=Surfacewater
GW=Groundwater
DW=Drinkingwater
S=Soil/Sludge
O=Oil
E=Extract
X=Other

Container Key
P=plastic G=glass

Preservation

Station Identification	Sample Date	Sample Time	Analysis	Unpres	4°C	HNO3	H2SO4	HCl	Methanol	Other	Matrix	Container number/type		pH	Analytics Sample #
MW-5	5/26/06	13:00	RCRA 8 Metals			✓					GW	1	250ml		56637-1
MW-5	5/26/06	13:00	SVOC 8270		✓						GW	1	1L		
MW-5	5/26/06	13:00	VOC					✓			GW	4	40ml		
* MW-4	5/26/06	07:30	RCRA 8 Metals			✓					GW	1	250ml		-2
MW-4	5/26/06	07:30	SVOC 8270		✓						GW	1	1L		
MW-4	5/26/06	07:30	VOC					✓			GW	4	40ml		
~~MW-3~~	~~5/26/06~~	~~07:30~~													
~~MW-3~~	~~5/26/06~~	07:30													
MW-3															

Received By: Time: 15:00 Date: 5/26/06
Received By: Time: 1815 Date: 5/26/06
Received By: Time: 955 Date: 5/27/06

Relinquished By Sampler:
Relinquished By:
Relinquished By:

FAX RESULTS? YES NO
Fax # 207-774-6635

Turnaround Request
Standard ☐ Due Date
Priority ☐ Due Date

Comments / Instructions:
Metal Sample Filtered
SVOC checked for residual Cl 0 PPM w/ KI paper 5/27/06
*MW-1 sample labled on bottle MW-4 reported as MW-4 as per KK as per 5-27-06 HSK
All GW Filtered as per KK

Page 1 of 1

Chain Of Custody Form

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Phone (603) 436-5111
Fax (603) 430-2151

For Analytics Use Only Rev. 1, 10/1/02

Samples were:
1) Shipped or hand-delivered
2) Temp blank °C 8.1°C/4
3) Received in good condition (Y) or N
4) pH checked by: MGL
5) Labels checked by: MJG 5.30.06

Project#: 21873G.002 Proj. Name: Cape Old Town Mill

Company: Woodard &Curran

Contact: Karl Keszler

Address: 41 Hutchins Drive
Portland, ME 04102

Phone: (207)774-2112 PO# Quote #

Sampler (Signature):

Matrix Key:
WW=Wastewater
SW=Surfacewater
GW=Groundwater
DW=Drinkingwater
S=Soil/Sludge
O=Oil
E=Extract
X=Other

Container Key
P=plastic G=glass

Preservation: Unpres | 4°C | HNO3 | H2SO4 | HCL | Methanol | Other

Station Identification	Sample Date	Sample Time	Analysis	Unpres	4°C	HNO3	H2SO4	HCL	Methanol	Other	Matrix	Container number/type		pH	Analytics Sample #
SB-37	5/26/06	14:25	RCRA 8 Metals			✓					GW	1	500ml P		506037-3
SB-37		14:25	SVOC 8270									1	1L G		
SB-37		14:25	VOC					✓				4	40ml G		
✓ SB-22 (0-4)	5/25/06	1600	PCB	✓							S	1	8oz		-4
✓ SB-22 (0-4)	5/25/06	1600	metals	✓							S	1	4oz		
✓ SB-22 (0-4)	5/25/06	1600	SVOCs	✓							S	1	4oz		
✓ SB-22 (0-4)	5/25/06	1600	VOCs												
~~SB-37 (0-4)~~	~~5/25/06~~	~~1500~~													

Comments / Instructions: Metals filtered. (GW) SB-37 SVOC checked for residual Cl w/KI paper DOM 5/27/06

FAX RESULTS? YES NO
Fax # 207-774-6635

Turnaround Request
Standard ☐ Due Date
Priority ☒ Due Date

Received By: Time: 15:00 Date: 5/26/06
Received By: Time: 1812 Date: 5/26/06
Received By: Time: 0855 Date: 5/27/06
Relinquished By Sampler:
Relinquished By:
Relinquished By:

Katahdin ANALYTICAL SERVICES
340 County Road No. 5
P.O. Box 520
Westbrook, ME 04092
Tel: (207) 874-2400
Fax: (207) 775-4029

Analytics

CHAIN of CUSTODY

PLEASE BEAR DOWN AND PRINT LEGIBLY IN PEN

Page ____ of ____

Client: Woodard + Curran | Contact: Karl Kasper | Phone #: (207) 774-2112 | Fax #: (207) 774-6635

Address: 41 Hutchins Drive | City: Portland | State: ME | Zip Code: 04102

Purchase Order #: 218736.002 | Proj. Name / No. | Katahdin Quote #

Bill (if different than above) | Address

Sampler (Print / Sign): Josh Bowe | Copies To:

LAB USE ONLY | WORK ORDER #:

KATAHDIN PROJECT NUMBER ____

REMARKS: ____

SHIPPING INFO: ☐ FED EX ☐ UPS ☐ CLIENT

AIRBILL NO: ____

TEMP°C 4 ☐ TEMP BLANK ☐ INTACT ☐ NOT INTACT

ANALYSIS AND CONTAINER TYPE PRESERVATIVES

*	Sample Description	Date / Time coll'd	Matrix	No. of Cntrs.	Filt. ☐Y ☑N: 40ml 8260 B unpreserved	Filt. ☐Y ☑N: PCB unpreserved 4oz	Filt. ☐Y ☑N: Metals RCRA8 unpreserved 8oz	Filt. ☐Y ☑N: 40ml low soil H2O preserved	Filt. ☐Y ☐N: 40ml med soil MeOH preserved	Filt. ☐Y ☐N: SVOA (8270) unpreserved 2oz	Filt. ☐Y ☐N: SVOA (8270) unpreserved 4oz	Filt. ☐Y ☐N	Filt. ☐Y ☐N	Filt. ☐Y ☐N
	SB-1 (6-10')	5/25/8:00	GW	2	2							56637 ~~56640~~ DPM 5/27/06		-5
	SB-21 (4-8')	/11:50	Soil	1		1								-6
	SB-35 (0-6")	/7:00	Soil	1		1								-7
	SB-5 (4-6')	/13:20	Soil	1		1								-8
	SB-21 (0-4')	/11:40	Soil	6		1	1	2	1	1				-9
	SB-5 (0-4')	/13:15	Soil	6		1	1	2	1	1				-10
	SB-6 (0-4')	/14:00	Soil	5			1	2	1		1			-11
	SB-23 (0-4)'	/15:30	Soil	3		1	1				1			-12
	SB-24 (0-4)'	/14:10	Soil	1		1								-13
		/												
		/												
		/												
		/												
		/												
		/												
		/												

COMMENTS: temp @ 4 OC when received by AEL DPM 5/26/06

Labels checked AEL by MKS 5-30-06

Relinquished By: (Signature)	Date / Time	Received By: (Signature)	Relinquished By: (Signature)	Date / Time	Received By: (Signature)
Josh Bowe	5/25 16:16	[signature]	[signature]	5/27/06 8:55	[signature]
Relinquished By: (Signature)	Date / Time	Received By: (Signature)	Relinquished By: (Signature)	Date / Time	Received By: (Signature)

THE TERMS AND CONDITIONS ON THE REVERSE SIDE HEREOF SHALL GOVERN SERVICES, EXCEPT WHEN A SIGNED CONTRACTUAL AGREEMENT EXISTS.

3 of 10 '1
ORIGINAL

Chain Of Custody Form

analytics environmental laboratory LLC
195 Commerce Way Suite E
Portsmouth, NH 03801
Phone (603) 436-5111
Fax (603) 430-2151

For Analytics Use Only Rev. 1, 10/1/02

Samples were:
1) Shipped or hand-delivered
2) Temp blank °C 8.1°C/4
3) Received in good condition (Y) or N
4) pH checked by: MEL
5) Labels checked by: MG 5.30.06

Project#: 218736 Proj. Name:
Company: Woodard &Curran
Contact: KARI KASPER
Address: 41 Hutchins Drive
Portland, ME 04102
Phone: (207)774-2112 PO# Quote #
Sampler (Signature): Dorothy Taylor

Matrix Key:
WW=Wastewater
SW=Surfacewater
GW=Groundwater
DW=Drinkingwater
S=Soil/Sludge
O=Oil
E=Extract
X=Other

Container Key
P=plastic G=glass

Preservation: Unpres | 4° C | HNO3 | H2SO4 | HCL | Methanol | Other

Station Identification	Sample Date	Sample Time	Analysis	Matrix	Container number/type	pH	Analytics Sample #
SB-31 (0-4)	5/25/06	10:30	8260 VOC	S	2 6		506040 - 14
"	"	"	8270 SVOC	S	1		506037
"	"	"	RCRA 8 Metals	S	1		
SB-30 (0-4)	"	11:25	8260 VOC		2		-15
"	"	11:25	8270 SVOC		1		
"	"	11:25	RCRA 8 Metals		1		
" DT	"		DT				
SB-38 (0-4)	"	13:00	8260 VOC		2		-16
"	"	13:00	8270 SVOC		1		
"	"	13:00	RCRA 8 Metals		1		
SB-37 (0-4)	"	13:40	8260 VOCs		2		-1617
"	"	13:40	8270 SVOC		1		DOM 5/27/06

Comments / Instructions:

FAX RESULTS? YES NO
fax # 207-774-6635

Turnaround Request
Standard ☐ Priority ☐
Due Date Due Date

Received By:	Time:	Date:	Relinquished By Sampler:
	1811	5/25/06	Dorothy Taylor
	855	5/27/06	

Page 4 of 7

Chain Of Custody Form

analytics environmental laboratory LLC
195 Commerce Way Suite E
Portsmouth, NH 03801
Phone (603) 436-5111
Fax (603) 430-2151

For Analytics Use Only Rev. 1, 10/1/02

Samples were:
1) Shipped or hand-delivered
2) Temp blank °C 4
3) Received in good condition (Y) or N
4) pH checked by: MEL
5) Labels checked by: MJS 5.30.06

Project#: 218736 Proj. Name:
Company: Woodard &Curran
Contact: KARI KASPER
Address: 41 Hutchins Drive
Portland, ME 04102
Phone: (207)774-2112 PO# Quote #
Sampler (Signature): Daniella Taylor

Matrix Key:
WW=Wastewater
SW=Surfacewater
GW=Groundwater
DW=Drinkingwater
S=Soil/Sludge
O=Oil
E=Extract
X=Other

Container Key
P=plastic G=glass

Preservation: Unpres | 4° C | HNO_3 | H_2SO_4 | HCL | Methanol | Other

Station Identification	Sample Date	Sample Time	Analysis	Matrix	Container number/type		pH	Analytics Sample #
SB-37 (0-4)	5/25/06	13:40	RCRA 8 Metals	S	1	G		506648 -17
SB-15 (0-4)	5/25/06	0920	metALS	S	1			506637 -18
SB-15 (0-4)	5/25/06	0930	SVOCs	S	1			
SB-15 (0-4)	5/25/06	0930	VOCs LS	S	3			
~~SB-40 (0-4)~~	~~5/26/06~~	~~0800~~	~~metals~~	~~S~~	~~1~~			
~~"~~	~~"~~	~~0800~~	~~SVOCs~~	~~S~~	~~1~~			
~~"~~	~~"~~	~~0820~~	~~VOCs~~	~~S~~	~~1~~			
SB-41 (0-4)		0850	metals	S	2			-19
			SVOCs	S	1			
			VOCs.	S	1			
			PCB	S	1			

Comments / Instructions:

TAX RESULTS? YES NO
Fax # 207-774-6635
Turnaround Request
Standard ☐ Priority ☐
Due Date Due Date

Page 1 of 3

Received By:	Received By:	Received By:
Time: 1630	Time: 1813	Time: 855
Date: 5/25/06	Date: 5/26/06	Date: 5/27/06
Relinquished By Sampler: Daniella Taylor	Relinquished By:	Relinquished By:

Chain Of Custody Form

analytics environmental laboratory LLC

195 Commerce Way Suite E
Portsmouth, NH 03801
Phone (603) 436-5111
Fax (603) 430-2151

For Analytics Use Only Rev. 1, 10/1/02

Samples were:

1) Shipped or hand-delivered
2) Temp blank °C 8.1°C / 4°C
3) Received in good condition (Y) or N
4) pH checked by: MEL
5) Labels checked by: 5.30.06

Project#: Proj. Name:

Company: Woodard &Curran

Contact:

Address: 41 Hutchins Drive

Portland, ME 04102

Phone: (207)774-2112 PO# Quote #

Sampler (Signature):

Matrix Key:
WW=Wastewater
SW=Surfacewater
GW=Groundwater
DW=Drinkingwater
S=Soil/Sludge
O=Oil
E=Extract
X=Other

Container Key
P=plastic G=glass

Station Identification	Sample Date	Sample Time	Analysis	Preservation: Unpres	4° C	HNO3	H2SO4	HCL	Methanol	Other	Matrix	Container number/type		pH	Analytics Sample #
SB-15 (0-4)	5/26/06	1310	SVOCs	X							GW S	1	G		5606 38 -20
SB-15 (0/4)		1330	VOCs							X	GW S DI	4			5606 37
SB-15 (0-4)		1330	Metals	X							GW S	1			
MW-3		1225	SVOCs	X							GW S	1			-21
MW-3		1225	VOCs							X	GW S DI	3			
MW-3		1250	metals	X							GW S	1			
Pile 1		1400	SVOCs	X							S	1			-22
Pile 1		1400	metals								S	1			
Pile 2		1410	SVOCs	X								1			-23
Pile 2		1410	Metals	X								1			
SB-40 (1½-5½)		0810	VOCs	X				X				2			-24
TPLP Blank	5/26		Voc					X				2			-29 -29

Comments / Instructions:

Metal GW samples filtered

SB-15 + MW-3 SVOC checked for residual Cl w/ KI paper

FAX RESULTS? YES NO

Fax # 207-774-6635

Turnaround Request

Standard ☐ Due Date

Priority ☐ Due Date

Received By: Time: Date:

Received By: Time: 1814 Date: 5/26/06

Received By: Time: 855 Date: 5/31/06

Relinquished By Sampler: Date:

Relinquished By: Date:

Relinquished By: Date:

Chain Of Custody Form

analytics environmental laboratory LLC

195 Commerce Way Suite E
Portsmouth, NH 03801
Phone (603) 436-5111
Fax (603) 430-2151

Project#: 218736 Proj. Name:

Company: Woodard &Curran

Contact: Karl Kasper

Address: 41 Hutchins Drive

Portland, ME 04102

Phone: (207)774-2112 PO# Quote #

Sampler (Signature): Ronald Taylor

Matrix Key:
WW=Wastewater
SW=Surfacewater
GW=Groundwater
DW=Drinkingwater
S=Soil/Sludge
O=Oil
E=Extract
X=Other

Preservation

Container Key
P=plastic G=glass

For Analytics Use Only Rev. 1, 10/1/02

Samples were:
1) Shipped or hand-delivered
2) Temp blank °C 8.1°C/4°C
3) Received in good condition (Y) or N
4) pH checked by: N/A
5) Labels checked by: MJS 5.30.06

Station Identification	Sample Date	Sample Time	Analysis	Unpres	4°C	HNO_3	H_2SO_4	HCL	Methanol	Other	Matrix	Container number/type		pH	Analytics Sample #
SB-3BDUP	5/25	1305	VOA	NONE							W				506636-25
~~SB~~ MW-5	5/26	1110	VOA	NONE							GW				506637-26
SB-30	5/25	1115	VOA	NONE							W				-27
SB-15	5/25	0940	VOA	NONE							W				-28

Comments / Instructions:

AX RESULTS? YES NO

ax # 207-774-6635

Turnaround Request

andard ☐ Priority ☐

ue Date Due Date

Page 7 of 27

Received By:	Relinquished By Sampler:
Time: 1810	Time:
Date: 5/26/06	Date:
Received By:	Relinquished By:
Time: 0830	Time:
Date: 5/27/06	Date:
Received By:	Relinquished By:
Time:	Time:
Date:	Date:

ANALYTICS SAMPLE RECEIPT CHECKLIST

analytics environmental laboratory LLC

AEL LAB#: 56637 COOLER NUMBER: _____

CLIENT: Woodard & Curran P NUMBER OF COOLERS: 2

PROJECT: GP Old Town Mill DATE RECEIVED: 5/27/06

A: PRELIMINARY EXAMINATION: DATE COOLER OPENED: 5/27/06

1. Cooler received by(initials) DPM Date Received: 5/27/06
2. Did cooler come with a shipping slip? Y N

If YES, enter carrier name and airbill number here: hand-delivered

3. Were custody seals on the outside of cooler? Y N
How many & where: _____ Seal Date: _____ Seal Name: _____

4. Did the custody seals arrive unbroken and intact upon arrival? Y N

5. COC#: _____

6. Were Custody papers filled out properly (ink,signed, etc)? (Y) N

7. Were custody papers sealed in a plastic bag? (Y) N

8. Did you sign the COC in the appropriate place? (Y) N

9. Was the project identifiable from the COC papers? (Y) N

10. Was enough ice used to chill the cooler? (Y) N Temp. of cooler: 4

B. Log-In: Date samples were logged in: DPM By: 5/27/06

11. Type of packing in cooler(bubble wrap, popcorn) (Y) N

12. Were all bottles sealed in separate plastic bags? (Y) N

13. Did all bottles arrive unbroken and were labels in good condition? (Y) N

14. Were all bottle labels complete(ID,Date,time,etc.) (Y) N

15. Did all bottle labels agree with custody papers? (Y) N

16. Were the correct containers used for the tests indicated: (Y) N

17. Were samples received at the correct pH? (Y) N

18. Was sufficient amount of sample sent for the tests indicated? (Y) N

19. Were bubbles absent in VOA samples? (Y) N

If NO, List sample #'s: _____

20. Laboratory labeling verified by (initials): Date: MJG 5-31-06

Katahdin
ANALYTICAL SERVICES

nelac
Cert. No. E87604

June 7, 2006

Ms. Erika Lloyd
Woodard & Curran
41 Hutchins Drive
Portland,ME 04102

RE: Katahdin Lab Number: WW2637
Project ID: Old Town
Project Manager: Mrs. Andrea Colby
Sample Receipt Date(s): May 31, 2006

Dear Ms. Lloyd:

Please find enclosed the following information:

* Report of Analysis (Analytical and/or Field)
* Quality Control Data Summary
* Chain of Custody (COC)
* Login Report

A copy of the Chain of Custody is included in the paginated report. The original COC is attached as an addendum to this report.

Should you have any questions or comments concerning this Report of Analysis, please do not hesitate to contact the project manager listed above. This cover letter is an integral part of the ROA.

We certify that the test results provided in this report meet all the requirements of the NELAC standards unless otherwise noted in an attached technical narrative or in the Report of Analysis.

We appreciate your continued use of our laboratory and look forward to working with you in the future. The following signature indicates technical review and acceptance of the data.

Sincerely,

KATAHDIN ANALYTICAL SERVICES

Leslie Dimond
Authorized Signature

6.7.06
Date

TECHNICAL NARRATIVE

Organics Laboratory

The samples of work order WW2637 were analyzed in accordance with "Test Methods for Evaluating Solid Wastes: Physical/Chemical Methods." SW-846. 2nd edition, 1982 (revised 1984), 3rd edition, 1986, and Updates I, II, IIA, and III 1996, Office of Solid Waste and Emergency Response, U.S. EPA, for the specific methods listed below or on the Report of Analysis. Some manual integrations may have been performed due to split peaks and/or corrected baselines. All have been flagged with an "M" (software-generated) on the pertinent quantitation reports.

8260B Analysis

Samples WW2637-1 and 11 were used for the matrix spike (MS) and matrix spike duplicate (MSD), as per client request.

The reported percent recovery acceptance limits for the Laboratory Control Samples (LCSs) are statistically derived for the full list of spiked compounds. The recoveries of the spiked analytes in the LCS, Matrix Spike (MS) and Matrix Spike Duplicate (MSD) are compared to these acceptance limits. Katahdin standard operating procedure is not to take corrective action until greater than ten percent of the spiked analytes (with the exception of the common laboratory contaminants acetone and methylene chloride) in the LCS are outside of the QC limits. If the associated MS/MSD has greater than ten percent of the spiked analytes outside of the QC limits, no corrective action is taken, as long the LCS is acceptable.

One or more target analytes may have been detected above the MDL in some method blanks. Any of these analytes that were also detected in any of the associated samples were flagged with a "B" qualifier indicating that the analyte was detected in the method blank analyzed and/or extracted concurrently with the sample.

The initial analysis of sample WW2637-10 had a low recovery for one surrogate that was outside of the laboratory established acceptance limits and had responses for four internal standards that resulted in %D's which were outside the laboratory acceptance limit of -50% to +100% of the response of the internal standard of the daily calibration verification standard. The reanalysis of this sample resulted in one low surrogate recovery and two internal standard responses, indicating a likely matrix effect. Therefore, only the first analysis is reported.

The initial analysis of sample WW2637-11 had one low surrogate recovery, which was outside of the laboratory established acceptance limits. The MS/MSD pair of this sample, WG28636-3 and WG28636-4, also had two low surrogate recoveries, indicating a likely matrix effect. Therefore, the sample and MS/MSD were not reanalyzed.

The initial analysis of sample WW2637-14 had one low surrogate recovery, which was outside of the laboratory established acceptance limits. The reanalysis of this sample had two low surrogate recoveries and had a low response for one internal standard that resulted in a %D which was outside the laboratory acceptance limit of -50% to +100% of the response of the internal standard of the daily calibration verification standard, indicating a likely matrix effect. Therefore, only the first analysis is reported.

8082 Analysis

The LCSD WG28624-3 had a high recovery for Aroclor 1260, which was outside of the laboratory established acceptance limits. Since the LCS WG28624-2, had an acceptable recovery for Aroclor 1260, the associated samples were not reextracted.

8270C Analysis

The reported percent recovery acceptance limits for the Laboratory Control Samples (LCSs) are statistically derived for the full list of spiked compounds. The recoveries of the spiked analytes in the LCS, Matrix Spike (MS) and Matrix Spike Duplicate (MSD) are compared to these acceptance limits. Katahdin standard operating procedure is not to take corrective action until greater than ten percent of the spiked analytes (with the exception of the common laboratory phthalate ester contaminants) in the LCS are outside of the QC limits. If the associated MS/MSD have greater than ten percent of the spiked analytes outside of the QC limits, no corrective action is taken, as long the LCS is acceptable.

Samples WW2637-1 and 11 were used for the matrix spike (MS) and matrix spike duplicate (MSD), as per client request.

The laboratory method blank WG28661-1 and samples WW2637-2 and 3 had low recoveries for one or two surrogates, which were outside the laboratory established acceptance limits. Since there were no second aliquots to reextract, no additional corrective action was taken.

Sample WW2637-11 had a low response for one internal standard, which was outside the laboratory acceptance limit of -50% to 100% of the response of the internal standard of the daily calibration verification standard. Based on the sample chromatogram, this deviation is likely attributable to a matrix effect. Therefore, the sample was not reanalyzed.

The laboratory method blank WG28761-1 had a high recovery for one surrogate, which was outside the laboratory established acceptance limits. Since a high recovery would indicate a high bias and there were no target analytes detected above the PQL, the associated sample was not reextracted.

The matrix spike sample (MS) WG28761-3 and the matrix spike duplicate sample (MSD) WG28761-4 had low responses for two internal standards, which were outside the laboratory acceptance limit of -50% to 100% of the response of the internal standard of the daily continuing calibration standard. Based on the sample chromatograms, the deviations are likely attributable to a matrix effect. Therefore, the samples were not reanalyzed.

There were no other protocol deviations or observations noted by the organics laboratory staff.

DATA QUALIFIERS

U	Indicates the compound was analyzed for but not detected above the laboratory Practical Quantitation Limit.
*	Compound recovery outside of quality control limits.
D	Indicates the result was obtained from analysis of a diluted sample. Surrogate recoveries may not be calculable.
E	Estimated value. This flag identifies compounds whose concentrations exceed the upper level of the calibration range of the instrument for that specific analysis.
J	Estimated value. The analyte was detected in the sample at a concentration less than the laboratory Practical Quantitation Limit (PQL), but above the Method Detection Limit (MDL).
B	Organics- Indicates the analyte was detected in the laboratory method blank analyzed concurrently with the sample. Metals- Indicates the analyte was detected in the sample at a concentration greater than the instrument detection limit, but less than the laboratory's Practical Quantitation Level.
N	Presumptive evidence of a compound based on a mass spectral library search.
A	Indicates that a tentatively identified compound is a suspected aldol-condensation product.
P	Used for Pesticide/Aroclor analyte when there is a greater than 25% difference for detected concentrations between the two GC columns.
MCL	Maximum Contaminant Level
NL	No limit
NFL	No Free Liquid Present
FLP	Free Liquid Present

KATAHDIN ANALYTICAL SERVICES
Report of Analytical Results

Client: Woodard & Curran
Project: Old Town
PO No:
Sample Date: 05/30/06
Received Date: 05/31/06
Extraction Date:
Analysis Date: 31-MAY-2006 15:37
Report Date: 06/02/2006
Matrix: WATER
% Solids: NA

Lab ID: WW2637-1
Client ID: SB-25
SDG: WW2637
Extracted by:
Extraction Method: SW846 5030
Analyst: SKT
Analysis Method: SW846 8260B
Lab Prep Batch: WG28646
Units: ug/l

Compound	Flags	Results	DF	PQL	Adj.PQL
Dichlorodifluoromethane	U	10	1.0	10	10
Chloromethane	U	10	1.0	10	10
Vinyl chloride	U	10	1.0	10	10
Bromomethane	U	10	1.0	10	10
Chloroethane	U	10	1.0	10	10
Trichlorofluoromethane	U	10	1.0	10	10
1,1-Dichloroethene	U	5	1.0	5	5
Methylene Chloride	U	5	1.0	5	5
trans-1,2-Dichloroethene	U	5	1.0	5	5
1,1-Dichloroethane	U	5	1.0	5	5
cis-1,2-Dichloroethene	U	5	1.0	5	5
1,2-Dichloroethylene (total)	U	10	1.0	10	10
2,2-Dichloropropane	U	5	1.0	5	5
Chloroform	U	5	1.0	5	5
Bromochloromethane	U	5	1.0	5	5
1,1,1-Trichloroethane	U	5	1.0	5	5
1,2-Dichloroethane	U	5	1.0	5	5
1,1-Dichloropropene	U	5	1.0	5	5
Carbon Tetrachloride	U	5	1.0	5	5
Benzene	U	5	1.0	5	5
1,2-Dichloropropane	U	5	1.0	5	5
Trichloroethene	U	5	1.0	5	5
Dibromomethane	U	5	1.0	5	5
Bromodichloromethane	U	5	1.0	5	5
cis-1,3-dichloropropene	U	5	1.0	5	5
Toluene	U	5	1.0	5	5
trans-1,3-Dichloropropene	U	5	1.0	5	5
1,1,2-Trichloroethane	U	5	1.0	5	5
1,3-Dichloropropane	U	5	1.0	5	5
Dibromochloromethane	U	5	1.0	5	5
Tetrachloroethene	U	5	1.0	5	5
1,2-Dibromoethane	U	5	1.0	5	5
Chlorobenzene	U	5	1.0	5	5
1,1,1,2-Tetrachloroethane	U	5	1.0	5	5
Ethylbenzene	U	5	1.0	5	5
Bromoform	U	5	1.0	5	5
Styrene	U	5	1.0	5	5
1,1,2,2-Tetrachloroethane	U	5	1.0	5	5
1,2,3-Trichloropropane	U	5	1.0	5	5
Isopropylbenzene	U	5	1.0	5	5
Bromobenzene	U	5	1.0	5	5
2-Chlorotoluene	U	5	1.0	5	5
N-Propylbenzene	U	5	1.0	5	5

KATAHDIN ANALYTICAL SERVICES
Report of Analytical Results

Client: Woodard & Curran
Project: Old Town
PO No:
Sample Date: 05/30/06
Received Date: 05/31/06
Extraction Date:
Analysis Date: 31-MAY-2006 15:37
Report Date: 06/02/2006
Matrix: WATER
% Solids: NA

Lab ID: WW2637-1
Client ID: SB-25
SDG: WW2637
Extracted by:
Extraction Method: SW846 5030
Analyst: SKT
Analysis Method: SW846 8260B
Lab Prep Batch: WG28646
Units: ug/l

Compound	Flags	Results	DF	PQL	Adj.PQL
4-Chlorotoluene	U	5	1.0	5	5
1,3,5-Trimethylbenzene	U	5	1.0	5	5
tert-Butylbenzene	U	5	1.0	5	5
1,2,4-Trichlorobenzene	U	5	1.0	5	5
sec-Butylbenzene	U	5	1.0	5	5
1,3-Dichlorobenzene	U	5	1.0	5	5
P-Isopropyltoluene	U	5	1.0	5	5
1,4-Dichlorobenzene	U	5	1.0	5	5
1,2-Dichlorobenzene	U	5	1.0	5	5
N-Butylbenzene	U	5	1.0	5	5
1,2-Dibromo-3-Chloropropane	U	5	1.0	5	5
1,2,4-Trimethylbenzene	U	5	1.0	5	5
Naphthalene	U	5	1.0	5	5
Hexachlorobutadiene	U	5	1.0	5	5
1,2,3-Trichlorobenzene	U	5	1.0	5	5
Methyl tert-butyl ether	U	10	1.0	10	10
Acetone	U	25	1.0	25	25
2-Butanone	U	25	1.0	25	25
4-methyl-2-pentanone	U	25	1.0	25	25
2-Hexanone	U	25	1.0	25	25
m+p-Xylenes	U	10	1.0	10	10
o-Xylene	U	5	1.0	5	5
Xylenes (total)	U	15	1.0	15	15
1,3,5-Trichlorobenzene	U	5	1.0	5	5
Vinyl Acetate	U	5	1.0	5	5
Carbon Disulfide	U	5	1.0	5	5
Diethyl Ether	U	5	1.0	5	5
Tetrahydrofuran	U	25	1.0	25	25
Dibromofluoromethane		91%			
1,2-Dichloroethane-D4		100%			
Toluene-D8		86%			
P-Bromofluorobenzene		87%			

KATAHDIN ANALYTICAL SERVICES
Report of Analytical Results

Client: Woodard & Curran
Project: Old Town
PO No:
Sample Date: 05/30/06
Received Date: 05/31/06
Extraction Date: 05/31/06
Analysis Date: 01-JUN-2006 15:17
Report Date: 06/06/2006
Matrix: WATER
% Solids: NA

Lab ID: WW2637-1
Client ID: SB-25
SDG: WW2637
Extracted by: GN
Extraction Method: SW846 3510
Analyst: LRS
Analysis Method: SW846 8270C
Lab Prep Batch: WG28661
Units: ug/L

Compound	Flags	Results	DF	PQL	Adj.PQL
Phenol	U	10	1.0	10	10
Bis(2-Chloroethyl)ether	U	10	1.0	10	10
2-Chlorophenol	U	10	1.0	10	10
1,3-Dichlorobenzene	U	10	1.0	10	10
1,4-Dichlorobenzene	U	10	1.0	10	10
1,2-Dichlorobenzene	U	10	1.0	10	10
2-Methylphenol	U	10	1.0	10	10
2,2'-Oxybis(1-Chloropropane)	U	10	1.0	10	10
N-Nitroso-di-n-propylamine	U	10	1.0	10	10
3&4-Methylphenol	U	10	1.0	10	10
Hexachloroethane	U	10	1.0	10	10
Nitrobenzene	U	10	1.0	10	10
Isophorone	U	10	1.0	10	10
2-Nitrophenol	U	10	1.0	10	10
2,4-Dimethylphenol	U	10	1.0	10	10
Bis(2-Chloroethoxy)methane	U	10	1.0	10	10
2,4-Dichlorophenol	U	10	1.0	10	10
1,2,4-Trichlorobenzene	U	10	1.0	10	10
Naphthalene	U	10	1.0	10	10
4-Chloroaniline	U	10	1.0	10	10
Hexachlorobutadiene	U	10	1.0	10	10
4-Chloro-3-Methylphenol	U	10	1.0	10	10
2-Methylnaphthalene	U	10	1.0	10	10
Hexachlorocyclopentadiene	U	10	1.0	10	10
2,4,6-Trichlorophenol	U	10	1.0	10	10
2,4,5-Trichlorophenol	U	25	1.0	25	25
2-Chloronaphthalene	U	10	1.0	10	10
2-Nitroaniline	U	25	1.0	25	25
Dimethyl Phthalate	U	10	1.0	10	10
2,6-Dinitrotoluene	U	10	1.0	10	10
Acenaphthylene	U	10	1.0	10	10
3-Nitroaniline	U	25	1.0	25	25
Acenaphthene	U	10	1.0	10	10
2,4-Dinitrophenol	U	25	1.0	25	25
Dibenzofuran	U	10	1.0	10	10
4-Nitrophenol	U	25	1.0	25	25
2,4-Dinitrotoluene	U	10	1.0	10	10
Diethylphthalate	U	10	1.0	10	10
Fluorene	U	10	1.0	10	10
4-Chlorophenyl-phenylether	U	10	1.0	10	10
4-Nitroaniline	U	25	1.0	25	25
4,6-Dinitro-2-Methylphenol	U	25	1.0	25	25
N-Nitrosodiphenylamine	U	10	1.0	10	10

KATAHDIN ANALYTICAL SERVICES
Report of Analytical Results

Client: Woodard & Curran
Project: Old Town
PO No:
Sample Date: 05/30/06
Received Date: 05/31/06
Extraction Date: 05/31/06
Analysis Date: 01-JUN-2006 15:17
Report Date: 06/06/2006
Matrix: WATER
% Solids: NA

Lab ID: WW2637-1
Client ID: SB-25
SDG: WW2637
Extracted by: GN
Extraction Method: SW846 3510
Analyst: LRS
Analysis Method: SW846 8270C
Lab Prep Batch: WG28661
Units: ug/L

Compound	Flags	Results	DF	PQL	Adj.PQL
4-Bromophenyl-phenylether	U	10	1.0	10	10
Hexachlorobenzene	U	10	1.0	10	10
Pentachlorophenol	U	25	1.0	25	25
Phenanthrene	U	10	1.0	10	10
Anthracene	U	10	1.0	10	10
Carbazole	U	10	1.0	10	10
Di-n-butylphthalate	U	10	1.0	10	10
Fluoranthene	U	10	1.0	10	10
Pyrene	U	10	1.0	10	10
Butylbenzylphthalate	U	10	1.0	10	10
Benzo(a)anthracene	U	10	1.0	10	10
3,3'-Dichlorobenzidine	U	10	1.0	10	10
Chrysene	U	10	1.0	10	10
bis(2-Ethylhexyl)phthalate	U	10	1.0	10	10
Di-n-octylphthalate	U	10	1.0	10	10
Benzo(b)fluoranthene	U	10	1.0	10	10
Benzo(k)fluoranthene	U	10	1.0	10	10
Benzo(a)pyrene	U	10	1.0	10	10
Indeno(1,2,3-cd)pyrene	U	10	1.0	10	10
Dibenzo(a,h)anthracene	U	10	1.0	10	10
Benzo(g,h,i)perylene	U	10	1.0	10	10
2-Fluorophenol		37%			
Phenol-D6		34%			
Nitrobenzene-D5		62%			
2-Fluorobiphenyl		65%			
2,4,6-Tribromophenol		58%			
Terphenyl-D14		82%			

Katahdin
ANALYTICAL SERVICES

REPORT OF ANALYTICAL RESULTS

Client: Erika Lloyd
Woodard & Curran
41 Hutchins Drive
Portland, ME 04102

Lab Sample ID: WW2637-001
Report Date: 6/7/2006
PO No.:
Project: Old Town

Sample Description	Matrix	Filtered	Date Sampled	Date Received
SB-25	AQ	No(Total)	05/30/2006	05/31/2006

Parameter	Result	Units	Adjusted PQL	Dilution Factor	PQL	Analytical Method	Analysis Date	By	Prep Method	Prepped Date	By	QC	Notes
ARSENIC	U 0.008	mg/L	0.008	1	0.008	SW846 6010	6/1/06	SJC	SW846 3010	5/31/06	MW	WE31ICW1	
BARIUM	0.0774	mg/L	0.0050	1	0.005	SW846 6010	5/31/06	SJC	SW846 3010	5/31/06	MW	WE31ICW1	
CADMIUM	U 0.0100	mg/L	0.0100	1	0.01	SW846 6010	5/31/06	SJC	SW846 3010	5/31/06	MW	WE31ICW1	
CHROMIUM	U 0.0030	mg/L	0.0030	5	0.0006	SW846 6020	6/2/06	EAM	SW846 3010	5/31/06	MW	WE31ICW1	1
LEAD	U 0.0010	mg/L	0.0010	5	0.0002	SW846 6020	6/2/06	EAM	SW846 3010	5/31/06	MW	WE31ICW1	1
MERCURY	U 0.20	ug/L	0.20	1	0.2	SW846 7470	6/1/06	MW	SW846 7470	6/1/06	MW	WF01HGW0	
SELENIUM	U 0.010	mg/L	0.010	1	0.01	SW846 6010	5/31/06	SJC	SW846 3010	5/31/06	MW	WE31ICW1	
SILVER	U 0.015	mg/L	0.015	1	0.015	SW846 6010	5/31/06	SJC	SW846 3010	5/31/06	MW	WE31ICW1	

1 The laboratory's Practical Quantitation Level could not be achieved for this parameter due to sample composition, matrix effects,sample volume, or quantity used for analysis.

KATAHDIN ANALYTICAL SERVICES
Report of Analytical Results

Client: Woodard & Curran
Project: Old Town
PO No:
Sample Date: 05/30/06
Received Date: 05/31/06
Extraction Date:
Analysis Date: 31-MAY-2006 16:08
Report Date: 06/02/2006
Matrix: WATER
% Solids: NA

Lab ID: WW2637-2
Client ID: MW-1
SDG: WW2637
Extracted by:
Extraction Method: SW846 5030
Analyst: SKT
Analysis Method: SW846 8260B
Lab Prep Batch: WG28646
Units: ug/l

Compound	Flags	Results	DF	PQL	Adj.PQL
Dichlorodifluoromethane	U	10	1.0	10	10
Chloromethane	U	10	1.0	10	10
Vinyl chloride	U	10	1.0	10	10
Bromomethane	U	10	1.0	10	10
Chloroethane	U	10	1.0	10	10
Trichlorofluoromethane	U	10	1.0	10	10
1,1-Dichloroethene	U	5	1.0	5	5
Methylene Chloride	U	5	1.0	5	5
trans-1,2-Dichloroethene	U	5	1.0	5	5
1,1-Dichloroethane	U	5	1.0	5	5
cis-1,2-Dichloroethene	U	5	1.0	5	5
1,2-Dichloroethylene (total)	U	10	1.0	10	10
2,2-Dichloropropane	U	5	1.0	5	5
Chloroform	U	5	1.0	5	5
Bromochloromethane	U	5	1.0	5	5
1,1,1-Trichloroethane	U	5	1.0	5	5
1,2-Dichloroethane	U	5	1.0	5	5
1,1-Dichloropropene	U	5	1.0	5	5
Carbon Tetrachloride	U	5	1.0	5	5
Benzene	U	5	1.0	5	5
1,2-Dichloropropane	U	5	1.0	5	5
Trichloroethene	U	5	1.0	5	5
Dibromomethane	U	5	1.0	5	5
Bromodichloromethane	U	5	1.0	5	5
cis-1,3-dichloropropene	U	5	1.0	5	5
Toluene	U	5	1.0	5	5
trans-1,3-Dichloropropene	U	5	1.0	5	5
1,1,2-Trichloroethane	U	5	1.0	5	5
1,3-Dichloropropane	U	5	1.0	5	5
Dibromochloromethane	U	5	1.0	5	5
Tetrachloroethene	U	5	1.0	5	5
1,2-Dibromoethane	U	5	1.0	5	5
Chlorobenzene	U	5	1.0	5	5
1,1,1,2-Tetrachloroethane	U	5	1.0	5	5
Ethylbenzene	U	5	1.0	5	5
Bromoform	U	5	1.0	5	5
Styrene	U	5	1.0	5	5
1,1,2,2-Tetrachloroethane	U	5	1.0	5	5
1,2,3-Trichloropropane	U	5	1.0	5	5
Isopropylbenzene	U	5	1.0	5	5
Bromobenzene	U	5	1.0	5	5
2-Chlorotoluene	U	5	1.0	5	5
N-Propylbenzene	U	5	1.0	5	5

KATAHDIN ANALYTICAL SERVICES
Report of Analytical Results

Client: Woodard & Curran
Project: Old Town
PO No:
Sample Date: 05/30/06
Received Date: 05/31/06
Extraction Date:
Analysis Date: 31-MAY-2006 16:08
Report Date: 06/02/2006
Matrix: WATER
% Solids: NA

Lab ID: WW2637-2
Client ID: MW-1
SDG: WW2637
Extracted by:
Extraction Method: SW846 5030
Analyst: SKT
Analysis Method: SW846 8260B
Lab Prep Batch: WG28646
Units: ug/l

Compound	Flags	Results	DF	PQL	Adj.PQL
4-Chlorotoluene	U	5	1.0	5	5
1,3,5-Trimethylbenzene	U	5	1.0	5	5
tert-Butylbenzene	U	5	1.0	5	5
1,2,4-Trichlorobenzene	U	5	1.0	5	5
sec-Butylbenzene	U	5	1.0	5	5
1,3-Dichlorobenzene	U	5	1.0	5	5
P-Isopropyltoluene	U	5	1.0	5	5
1,4-Dichlorobenzene	U	5	1.0	5	5
1,2-Dichlorobenzene	U	5	1.0	5	5
N-Butylbenzene	U	5	1.0	5	5
1,2-Dibromo-3-Chloropropane	U	5	1.0	5	5
1,2,4-Trimethylbenzene	U	5	1.0	5	5
Naphthalene	U	5	1.0	5	5
Hexachlorobutadiene	U	5	1.0	5	5
1,2,3-Trichlorobenzene	U	5	1.0	5	5
Methyl tert-butyl ether	U	10	1.0	10	10
Acetone	U	25	1.0	25	25
2-Butanone	U	25	1.0	25	25
4-methyl-2-pentanone	U	25	1.0	25	25
2-Hexanone	U	25	1.0	25	25
m+p-Xylenes	U	10	1.0	10	10
o-Xylene	U	5	1.0	5	5
Xylenes (total)	U	15	1.0	15	15
1,3,5-Trichlorobenzene	U	5	1.0	5	5
Vinyl Acetate	U	5	1.0	5	5
Carbon Disulfide	U	5	1.0	5	5
Diethyl Ether	U	5	1.0	5	5
Tetrahydrofuran	U	25	1.0	25	25
Dibromofluoromethane		92%			
1,2-Dichloroethane-D4		105%			
Toluene-D8		88%			
P-Bromofluorobenzene		88%			

KATAHDIN ANALYTICAL SERVICES
Report of Analytical Results

Client: Woodard & Curran
Project: Old Town
PO No:
Sample Date: 05/30/06
Received Date: 05/31/06
Extraction Date: 05/31/06
Analysis Date: 01-JUN-2006 16:04
Report Date: 06/06/2006
Matrix: WATER
% Solids: NA

Lab ID: WW2637-2
Client ID: MW-1
SDG: WW2637
Extracted by: GN
Extraction Method: SW846 3510
Analyst: LRS
Analysis Method: SW846 8270C
Lab Prep Batch: WG28661
Units: ug/L

Compound	Flags	Results	DF	PQL	Adj.PQL
Phenol	U	10	1.0	10	10
Bis(2-Chloroethyl)ether	U	10	1.0	10	10
2-Chlorophenol	U	10	1.0	10	10
1,3-Dichlorobenzene	U	10	1.0	10	10
1,4-Dichlorobenzene	U	10	1.0	10	10
1,2-Dichlorobenzene	U	10	1.0	10	10
2-Methylphenol	U	10	1.0	10	10
2,2'-Oxybis(1-Chloropropane)	U	10	1.0	10	10
N-Nitroso-di-n-propylamine	U	10	1.0	10	10
3&4-Methylphenol	U	10	1.0	10	10
Hexachloroethane	U	10	1.0	10	10
Nitrobenzene	U	10	1.0	10	10
Isophorone	U	10	1.0	10	10
2-Nitrophenol	U	10	1.0	10	10
2,4-Dimethylphenol	U	10	1.0	10	10
Bis(2-Chloroethoxy)methane	U	10	1.0	10	10
2,4-Dichlorophenol	U	10	1.0	10	10
1,2,4-Trichlorobenzene	U	10	1.0	10	10
Naphthalene	U	10	1.0	10	10
4-Chloroaniline	U	10	1.0	10	10
Hexachlorobutadiene	U	10	1.0	10	10
4-Chloro-3-Methylphenol	U	10	1.0	10	10
2-Methylnaphthalene	U	10	1.0	10	10
Hexachlorocyclopentadiene	U	10	1.0	10	10
2,4,6-Trichlorophenol	U	10	1.0	10	10
2,4,5-Trichlorophenol	U	25	1.0	25	25
2-Chloronaphthalene	U	10	1.0	10	10
2-Nitroaniline	U	25	1.0	25	25
Dimethyl Phthalate	U	10	1.0	10	10
2,6-Dinitrotoluene	U	10	1.0	10	10
Acenaphthylene	U	10	1.0	10	10
3-Nitroaniline	U	25	1.0	25	25
Acenaphthene	U	10	1.0	10	10
2,4-Dinitrophenol	U	25	1.0	25	25
Dibenzofuran	U	10	1.0	10	10
4-Nitrophenol	U	25	1.0	25	25
2,4-Dinitrotoluene	U	10	1.0	10	10
Diethylphthalate	U	10	1.0	10	10
Fluorene	U	10	1.0	10	10
4-Chlorophenyl-phenylether	U	10	1.0	10	10
4-Nitroaniline	U	25	1.0	25	25
4,6-Dinitro-2-Methylphenol	U	25	1.0	25	25
N-Nitrosodiphenylamine	U	10	1.0	10	10

KATAHDIN ANALYTICAL SERVICES
Report of Analytical Results

Client: Woodard & Curran
Project: Old Town
PO No:
Sample Date: 05/30/06
Received Date: 05/31/06
Extraction Date: 05/31/06
Analysis Date: 01-JUN-2006 16:04
Report Date: 06/06/2006
Matrix: WATER
% Solids: NA

Lab ID: WW2637-2
Client ID: MW-1
SDG: WW2637
Extracted by: GN
Extraction Method: SW846 3510
Analyst: LRS
Analysis Method: SW846 8270C
Lab Prep Batch: WG28661
Units: ug/L

Compound	Flags	Results	DF	PQL	Adj.PQL
4-Bromophenyl-phenylether	U	10	1.0	10	10
Hexachlorobenzene	U	10	1.0	10	10
Pentachlorophenol	U	25	1.0	25	25
Phenanthrene	U	10	1.0	10	10
Anthracene	U	10	1.0	10	10
Carbazole	U	10	1.0	10	10
Di-n-butylphthalate	U	10	1.0	10	10
Fluoranthene	U	10	1.0	10	10
Pyrene	U	10	1.0	10	10
Butylbenzylphthalate	U	10	1.0	10	10
Benzo(a)anthracene	U	10	1.0	10	10
3,3'-Dichlorobenzidine	U	10	1.0	10	10
Chrysene	U	10	1.0	10	10
bis(2-Ethylhexyl)phthalate	U	10	1.0	10	10
Di-n-octylphthalate	U	10	1.0	10	10
Benzo(b)fluoranthene	U	10	1.0	10	10
Benzo(k)fluoranthene	U	10	1.0	10	10
Benzo(a)pyrene	U	10	1.0	10	10
Indeno(1,2,3-cd)pyrene	U	10	1.0	10	10
Dibenzo(a,h)anthracene	U	10	1.0	10	10
Benzo(g,h,i)perylene	U	10	1.0	10	10
2-Fluorophenol		24%			
Phenol-D6		24%			
Nitrobenzene-D5		69%			
2-Fluorobiphenyl		67%			
2,4,6-Tribromophenol		* 41%			
Terphenyl-D14		80%			

Katahdin
ANALYTICAL SERVICES

REPORT OF ANALYTICAL RESULTS

Client: Erika Lloyd
Woodard & Curran
41 Hutchins Drive
Portland, ME 04102

Lab Sample ID: WW2637-002
Report Date: 6/7/2006
PO No.:
Project: Old Town

Sample Description	Matrix	Filtered	Date Sampled	Date Received
MW-1	AQ	No(Total)	05/30/2006	05/31/2006

Parameter	Result	Units	Adjusted PQL	Dilution Factor	PQL	Analytical Method	Analysis Date	By	Prep Method	Prepped Date	By	QC	Notes
ARSENIC	U 0.0050	mg/L	0.0050	5	0.001	SW846 6020	6/2/06	EAM	SW846 3010	5/31/06	MW	WE31ICW1	1
BARIUM	0.104	mg/L	0.0050	1	0.005	SW846 6010	5/31/06	SJC	SW846 3010	5/31/06	MW	WE31ICW1	
CADMIUM	U 0.0100	mg/L	0.0100	1	0.01	SW846 6010	5/31/06	SJC	SW846 3010	5/31/06	MW	WE31ICW1	
CHROMIUM	U 0.0030	mg/L	0.0030	5	0.0006	SW846 6020	6/2/06	EAM	SW846 3010	5/31/06	MW	WE31ICW1	1
LEAD	U 0.0010	mg/L	0.0010	5	0.0002	SW846 6020	6/2/06	EAM	SW846 3010	5/31/06	MW	WE31ICW1	1
MERCURY	U 0.20	ug/L	0.20	1	0.2	SW846 7470	6/1/06	MW	SW846 7470	6/1/06	MW	WF01HGW0	
SELENIUM	U 0.010	mg/L	0.010	1	0.01	SW846 6010	5/31/06	SJC	SW846 3010	5/31/06	MW	WE31ICW1	
SILVER	U 0.015	mg/L	0.015	1	0.015	SW846 6010	5/31/06	SJC	SW846 3010	5/31/06	MW	WE31ICW1	

1 The laboratory's Practical Quantitation Level could not be achieved for this parameter due to sample composition, matrix effects,sample volume, or quantity used for analysis.

KATAHDIN ANALYTICAL SERVICES
Report of Analytical Results

Client: Woodard & Curran
Project: Old Town
PO No:
Sample Date: 05/30/06
Received Date: 05/31/06
Extraction Date:
Analysis Date: 31-MAY-2006 16:38
Report Date: 06/02/2006
Matrix: WATER
% Solids: NA

Lab ID: WW2637-3
Client ID: MW-2
SDG: WW2637
Extracted by:
Extraction Method: SW846 5030
Analyst: SKT
Analysis Method: SW846 8260B
Lab Prep Batch: WG28646
Units: ug/l

Compound	Flags	Results	DF	PQL	Adj.PQL
Dichlorodifluoromethane	U	10	1.0	10	10
Chloromethane	U	10	1.0	10	10
Vinyl chloride	U	10	1.0	10	10
Bromomethane	U	10	1.0	10	10
Chloroethane	U	10	1.0	10	10
Trichlorofluoromethane	U	10	1.0	10	10
1,1-Dichloroethene	U	5	1.0	5	5
Methylene Chloride	U	5	1.0	5	5
trans-1,2-Dichloroethene	U	5	1.0	5	5
1,1-Dichloroethane	U	5	1.0	5	5
cis-1,2-Dichloroethene	U	5	1.0	5	5
1,2-Dichloroethylene (total)	U	10	1.0	10	10
2,2-Dichloropropane	U	5	1.0	5	5
Chloroform	U	5	1.0	5	5
Bromochloromethane	U	5	1.0	5	5
1,1,1-Trichloroethane	U	5	1.0	5	5
1,2-Dichloroethane	U	5	1.0	5	5
1,1-Dichloropropene	U	5	1.0	5	5
Carbon Tetrachloride	U	5	1.0	5	5
Benzene	U	5	1.0	5	5
1,2-Dichloropropane	U	5	1.0	5	5
Trichloroethene	U	5	1.0	5	5
Dibromomethane	U	5	1.0	5	5
Bromodichloromethane	U	5	1.0	5	5
cis-1,3-dichloropropene	U	5	1.0	5	5
Toluene	U	5	1.0	5	5
trans-1,3-Dichloropropene	U	5	1.0	5	5
1,1,2-Trichloroethane	U	5	1.0	5	5
1,3-Dichloropropane	U	5	1.0	5	5
Dibromochloromethane	U	5	1.0	5	5
Tetrachloroethene	U	5	1.0	5	5
1,2-Dibromoethane	U	5	1.0	5	5
Chlorobenzene	U	5	1.0	5	5
1,1,1,2-Tetrachloroethane	U	5	1.0	5	5
Ethylbenzene	U	5	1.0	5	5
Bromoform	U	5	1.0	5	5
Styrene	U	5	1.0	5	5
1,1,2,2-Tetrachloroethane	U	5	1.0	5	5
1,2,3-Trichloropropane	U	5	1.0	5	5
Isopropylbenzene	U	5	1.0	5	5
Bromobenzene	U	5	1.0	5	5
2-Chlorotoluene	U	5	1.0	5	5
N-Propylbenzene	U	5	1.0	5	5

KATAHDIN ANALYTICAL SERVICES
Report of Analytical Results

Client: Woodard & Curran
Project: Old Town
PO No:
Sample Date: 05/30/06
Received Date: 05/31/06
Extraction Date:
Analysis Date: 31-MAY-2006 16:38
Report Date: 06/02/2006
Matrix: WATER
% Solids: NA

Lab ID: WW2637-3
Client ID: MW-2
SDG: WW2637
Extracted by:
Extraction Method: SW846 5030
Analyst: SKT
Analysis Method: SW846 8260B
Lab Prep Batch: WG28646
Units: ug/l

Compound	Flags	Results	DF	PQL	Adj.PQL
4-Chlorotoluene	U	5	1.0	5	5
1,3,5-Trimethylbenzene	U	5	1.0	5	5
tert-Butylbenzene	U	5	1.0	5	5
1,2,4-Trichlorobenzene	U	5	1.0	5	5
sec-Butylbenzene	U	5	1.0	5	5
1,3-Dichlorobenzene	U	5	1.0	5	5
P-Isopropyltoluene	U	5	1.0	5	5
1,4-Dichlorobenzene	U	5	1.0	5	5
1,2-Dichlorobenzene	U	5	1.0	5	5
N-Butylbenzene	U	5	1.0	5	5
1,2-Dibromo-3-Chloropropane	U	5	1.0	5	5
1,2,4-Trimethylbenzene	U	5	1.0	5	5
Naphthalene	U	5	1.0	5	5
Hexachlorobutadiene	U	5	1.0	5	5
1,2,3-Trichlorobenzene	U	5	1.0	5	5
Methyl tert-butyl ether	U	10	1.0	10	10
Acetone	U	25	1.0	25	25
2-Butanone	U	25	1.0	25	25
4-methyl-2-pentanone	U	25	1.0	25	25
2-Hexanone	U	25	1.0	25	25
m+p-Xylenes	U	10	1.0	10	10
o-Xylene	U	5	1.0	5	5
Xylenes (total)	U	15	1.0	15	15
1,3,5-Trichlorobenzene	U	5	1.0	5	5
Vinyl Acetate	U	5	1.0	5	5
Carbon Disulfide	U	5	1.0	5	5
Diethyl Ether	U	5	1.0	5	5
Tetrahydrofuran	U	25	1.0	25	25
Dibromofluoromethane		93%			
1,2-Dichloroethane-D4		102%			
Toluene-D8		86%			
P-Bromofluorobenzene		86%			

KATAHDIN ANALYTICAL SERVICES
Report of Analytical Results

Client: Woodard & Curran
Project: Old Town
PO No:
Sample Date: 05/30/06
Received Date: 05/31/06
Extraction Date: 05/31/06
Analysis Date: 01-JUN-2006 16:51
Report Date: 06/06/2006
Matrix: WATER
% Solids: NA

Lab ID: WW2637-3
Client ID: MW-2
SDG: WW2637
Extracted by: GN
Extraction Method: SW846 3510
Analyst: LRS
Analysis Method: SW846 8270C
Lab Prep Batch: WG28661
Units: ug/L

Compound	Flags	Results	DF	PQL	Adj.PQL
Phenol	U	10	1.0	10	10
Bis(2-Chloroethyl)ether	U	10	1.0	10	10
2-Chlorophenol	U	10	1.0	10	10
1,3-Dichlorobenzene	U	10	1.0	10	10
1,4-Dichlorobenzene	U	10	1.0	10	10
1,2-Dichlorobenzene	U	10	1.0	10	10
2-Methylphenol	U	10	1.0	10	10
2,2'-Oxybis(1-Chloropropane)	U	10	1.0	10	10
N-Nitroso-di-n-propylamine	U	10	1.0	10	10
3&4-Methylphenol	U	10	1.0	10	10
Hexachloroethane	U	10	1.0	10	10
Nitrobenzene	U	10	1.0	10	10
Isophorone	U	10	1.0	10	10
2-Nitrophenol	U	10	1.0	10	10
2,4-Dimethylphenol	U	10	1.0	10	10
Bis(2-Chloroethoxy)methane	U	10	1.0	10	10
2,4-Dichlorophenol	U	10	1.0	10	10
1,2,4-Trichlorobenzene	U	10	1.0	10	10
Naphthalene	U	10	1.0	10	10
4-Chloroaniline	U	10	1.0	10	10
Hexachlorobutadiene	U	10	1.0	10	10
4-Chloro-3-Methylphenol	U	10	1.0	10	10
2-Methylnaphthalene	U	10	1.0	10	10
Hexachlorocyclopentadiene	U	10	1.0	10	10
2,4,6-Trichlorophenol	U	10	1.0	10	10
2,4,5-Trichlorophenol	U	25	1.0	25	25
2-Chloronaphthalene	U	10	1.0	10	10
2-Nitroaniline	U	25	1.0	25	25
Dimethyl Phthalate	U	10	1.0	10	10
2,6-Dinitrotoluene	U	10	1.0	10	10
Acenaphthylene	U	10	1.0	10	10
3-Nitroaniline	U	25	1.0	25	25
Acenaphthene	U	10	1.0	10	10
2,4-Dinitrophenol	U	25	1.0	25	25
Dibenzofuran	U	10	1.0	10	10
4-Nitrophenol	U	25	1.0	25	25
2,4-Dinitrotoluene	U	10	1.0	10	10
Diethylphthalate	U	10	1.0	10	10
Fluorene	U	10	1.0	10	10
4-Chlorophenyl-phenylether	U	10	1.0	10	10
4-Nitroaniline	U	25	1.0	25	25
4,6-Dinitro-2-Methylphenol	U	25	1.0	25	25
N-Nitrosodiphenylamine	U	10	1.0	10	10

KATAHDIN ANALYTICAL SERVICES
Report of Analytical Results

Client: Woodard & Curran
Project: Old Town
PO No:
Sample Date: 05/30/06
Received Date: 05/31/06
Extraction Date: 05/31/06
Analysis Date: 01-JUN-2006 16:51
Report Date: 06/06/2006
Matrix: WATER
% Solids: NA

Lab ID: WW2637-3
Client ID: MW-2
SDG: WW2637
Extracted by: GN
Extraction Method: SW846 3510
Analyst: LRS
Analysis Method: SW846 8270C
Lab Prep Batch: WG28661
Units: ug/L

Compound	Flags	Results	DF	PQL	Adj.PQL
4-Bromophenyl-phenylether	U	10	1.0	10	10
Hexachlorobenzene	U	10	1.0	10	10
Pentachlorophenol	U	25	1.0	25	25
Phenanthrene	U	10	1.0	10	10
Anthracene	U	10	1.0	10	10
Carbazole	U	10	1.0	10	10
Di-n-butylphthalate	U	10	1.0	10	10
Fluoranthene	U	10	1.0	10	10
Pyrene	U	10	1.0	10	10
Butylbenzylphthalate	U	10	1.0	10	10
Benzo(a)anthracene	U	10	1.0	10	10
3,3'-Dichlorobenzidine	U	10	1.0	10	10
Chrysene	U	10	1.0	10	10
bis(2-Ethylhexyl)phthalate	U	10	1.0	10	10
Di-n-octylphthalate	U	10	1.0	10	10
Benzo(b)fluoranthene	U	10	1.0	10	10
Benzo(k)fluoranthene	U	10	1.0	10	10
Benzo(a)pyrene	U	10	1.0	10	10
Indeno(1,2,3-cd)pyrene	U	10	1.0	10	10
Dibenzo(a,h)anthracene	U	10	1.0	10	10
Benzo(g,h,i)perylene	U	10	1.0	10	10
2-Fluorophenol		34%			
Phenol-D6		29%			
Nitrobenzene-D5		* 48%			
2-Fluorobiphenyl		* 54%			
2,4,6-Tribromophenol		61%			
Terphenyl-D14		82%			

Katahdin
ANALYTICAL SERVICES

REPORT OF ANALYTICAL RESULTS

Client: Erika Lloyd
Woodard & Curran
41 Hutchins Drive
Portland, ME 04102

Lab Sample ID: WW2637-003
Report Date: 6/7/2006
PO No.:
Project: Old Town

Sample Description	Matrix	Filtered	Date Sampled	Date Received
MW-2	AQ	No(Total)	05/30/2006	05/31/2006

Parameter	Result	Units	Adjusted PQL	Dilution Factor	PQL	Analytical Method	Analysis Date	By	Prep Method	Prepped Date	By	QC	Notes
ARSENIC	U 0.0050	mg/L	0.0050	5	0.001	SW846 6020	6/2/06	EAM	SW846 3010	5/31/06	MW	WE31ICW1	1
BARIUM	0.0427	mg/L	0.0050	1	0.005	SW846 6010	5/31/06	SJC	SW846 3010	5/31/06	MW	WE31ICW1	
CADMIUM	U 0.0100	mg/L	0.0100	1	0.01	SW846 6010	5/31/06	SJC	SW846 3010	5/31/06	MW	WE31ICW1	
CHROMIUM	U 0.0030	mg/L	0.0030	5	0.0006	SW846 6020	6/2/06	EAM	SW846 3010	5/31/06	MW	WE31ICW1	1
LEAD	U 0.0010	mg/L	0.0010	5	0.0002	SW846 6020	6/2/06	EAM	SW846 3010	5/31/06	MW	WE31ICW1	1
MERCURY	U 0.20	ug/L	0.20	1	0.2	SW846 7470	6/1/06	MW	SW846 7470	6/1/06	MW	WF01HGW0	
SELENIUM	U 0.010	mg/L	0.010	1	0.01	SW846 6010	5/31/06	SJC	SW846 3010	5/31/06	MW	WE31ICW1	
SILVER	U 0.015	mg/L	0.015	1	0.015	SW846 6010	5/31/06	SJC	SW846 3010	5/31/06	MW	WE31ICW1	

1 The laboratory's Practical Quantitation Level could not be achieved for this parameter due to sample composition, matrix effects,sample volume, or quantity used for analysis.

KATAHDIN ANALYTICAL SERVICES
Report of Analytical Results

Client: Woodard & Curran
Project: Old Town
PO No:
Sample Date: 05/30/06
Received Date: 05/31/06
Extraction Date:
Analysis Date: 31-MAY-2006 17:09
Report Date: 06/02/2006
Matrix: WATER
% Solids: NA

Lab ID: WW2637-4
Client ID: SB-26
SDG: WW2637
Extracted by:
Extraction Method: SW846 5030
Analyst: SKT
Analysis Method: SW846 8260B
Lab Prep Batch: WG28646
Units: ug/l

Compound	Flags	Results	DF	PQL	Adj.PQL
Dichlorodifluoromethane	U	10	1.0	10	10
Chloromethane	U	10	1.0	10	10
Vinyl chloride	U	10	1.0	10	10
Bromomethane	U	10	1.0	10	10
Chloroethane	U	10	1.0	10	10
Trichlorofluoromethane	U	10	1.0	10	10
1,1-Dichloroethene	U	5	1.0	5	5
Methylene Chloride	U	5	1.0	5	5
trans-1,2-Dichloroethene	U	5	1.0	5	5
1,1-Dichloroethane	U	5	1.0	5	5
cis-1,2-Dichloroethene	U	5	1.0	5	5
1,2-Dichloroethylene (total)	U	10	1.0	10	10
2,2-Dichloropropane	U	5	1.0	5	5
Chloroform	U	5	1.0	5	5
Bromochloromethane	U	5	1.0	5	5
1,1,1-Trichloroethane	U	5	1.0	5	5
1,2-Dichloroethane	U	5	1.0	5	5
1,1-Dichloropropene	U	5	1.0	5	5
Carbon Tetrachloride	U	5	1.0	5	5
Benzene	U	5	1.0	5	5
1,2-Dichloropropane	U	5	1.0	5	5
Trichloroethene	U	5	1.0	5	5
Dibromomethane	U	5	1.0	5	5
Bromodichloromethane	U	5	1.0	5	5
cis-1,3-dichloropropene	U	5	1.0	5	5
Toluene	U	5	1.0	5	5
trans-1,3-Dichloropropene	U	5	1.0	5	5
1,1,2-Trichloroethane	U	5	1.0	5	5
1,3-Dichloropropane	U	5	1.0	5	5
Dibromochloromethane	U	5	1.0	5	5
Tetrachloroethene	U	5	1.0	5	5
1,2-Dibromoethane	U	5	1.0	5	5
Chlorobenzene	U	5	1.0	5	5
1,1,1,2-Tetrachloroethane	U	5	1.0	5	5
Ethylbenzene	U	5	1.0	5	5
Bromoform	U	5	1.0	5	5
Styrene	U	5	1.0	5	5
1,1,2,2-Tetrachloroethane	U	5	1.0	5	5
1,2,3-Trichloropropane	U	5	1.0	5	5
Isopropylbenzene	U	5	1.0	5	5
Bromobenzene	U	5	1.0	5	5
2-Chlorotoluene	U	5	1.0	5	5
N-Propylbenzene	U	5	1.0	5	5

KATAHDIN ANALYTICAL SERVICES
Report of Analytical Results

Client: Woodard & Curran
Project: Old Town
PO No:
Sample Date: 05/30/06
Received Date: 05/31/06
Extraction Date:
Analysis Date: 31-MAY-2006 17:09
Report Date: 06/02/2006
Matrix: WATER
% Solids: NA

Lab ID: WW2637-4
Client ID: SB-26
SDG: WW2637
Extracted by:
Extraction Method: SW846 5030
Analyst: SKT
Analysis Method: SW846 8260B
Lab Prep Batch: WG28646
Units: ug/l

Compound	Flags	Results	DF	PQL	Adj.PQL
4-Chlorotoluene	U	5	1.0	5	5
1,3,5-Trimethylbenzene	U	5	1.0	5	5
tert-Butylbenzene	U	5	1.0	5	5
1,2,4-Trichlorobenzene	U	5	1.0	5	5
sec-Butylbenzene	U	5	1.0	5	5
1,3-Dichlorobenzene	U	5	1.0	5	5
P-Isopropyltoluene	U	5	1.0	5	5
1,4-Dichlorobenzene	U	5	1.0	5	5
1,2-Dichlorobenzene	U	5	1.0	5	5
N-Butylbenzene	U	5	1.0	5	5
1,2-Dibromo-3-Chloropropane	U	5	1.0	5	5
1,2,4-Trimethylbenzene	U	5	1.0	5	5
Naphthalene	U	5	1.0	5	5
Hexachlorobutadiene	U	5	1.0	5	5
1,2,3-Trichlorobenzene	U	5	1.0	5	5
Methyl tert-butyl ether	U	10	1.0	10	10
Acetone	U	25	1.0	25	25
2-Butanone	U	25	1.0	25	25
4-methyl-2-pentanone	U	25	1.0	25	25
2-Hexanone	U	25	1.0	25	25
m+p-Xylenes	U	10	1.0	10	10
o-Xylene	U	5	1.0	5	5
Xylenes (total)	U	15	1.0	15	15
1,3,5-Trichlorobenzene	U	5	1.0	5	5
Vinyl Acetate	U	5	1.0	5	5
Carbon Disulfide	U	5	1.0	5	5
Diethyl Ether	U	5	1.0	5	5
Tetrahydrofuran	U	25	1.0	25	25
Dibromofluoromethane		91%			
1,2-Dichloroethane-D4		100%			
Toluene-D8		85%			
P-Bromofluorobenzene		85%			

KATAHDIN ANALYTICAL SERVICES
Report of Analytical Results

Client: Woodard & Curran
Project: Old Town
PO No:
Sample Date: 05/30/06
Received Date: 05/31/06
Extraction Date: 05/31/06
Analysis Date: 02-JUN-2006 08:15
Report Date: 06/06/2006
Matrix: WATER
% Solids: NA

Lab ID: WW2637-4
Client ID: SB-26
SDG: WW2637
Extracted by: GN
Extraction Method: SW846 3510
Analyst: LRS
Analysis Method: SW846 8270C
Lab Prep Batch: WG28661
Units: ug/L

Compound	Flags	Results	DF	PQL	Adj.PQL
Phenol	U	10	1.0	10	10
Bis(2-Chloroethyl)ether	U	10	1.0	10	10
2-Chlorophenol	U	10	1.0	10	10
1,3-Dichlorobenzene	U	10	1.0	10	10
1,4-Dichlorobenzene	U	10	1.0	10	10
1,2-Dichlorobenzene	U	10	1.0	10	10
2-Methylphenol	U	10	1.0	10	10
2,2'-Oxybis(1-Chloropropane)	U	10	1.0	10	10
N-Nitroso-di-n-propylamine	U	10	1.0	10	10
3&4-Methylphenol	U	10	1.0	10	10
Hexachloroethane	U	10	1.0	10	10
Nitrobenzene	U	10	1.0	10	10
Isophorone	U	10	1.0	10	10
2-Nitrophenol	U	10	1.0	10	10
2,4-Dimethylphenol	U	10	1.0	10	10
Bis(2-Chloroethoxy)methane	U	10	1.0	10	10
2,4-Dichlorophenol	U	10	1.0	10	10
1,2,4-Trichlorobenzene	U	10	1.0	10	10
Naphthalene	U	10	1.0	10	10
4-Chloroaniline	U	10	1.0	10	10
Hexachlorobutadiene	U	10	1.0	10	10
4-Chloro-3-Methylphenol	U	10	1.0	10	10
2-Methylnaphthalene	U	10	1.0	10	10
Hexachlorocyclopentadiene	U	10	1.0	10	10
2,4,6-Trichlorophenol	U	10	1.0	10	10
2,4,5-Trichlorophenol	U	25	1.0	25	25
2-Chloronaphthalene	U	10	1.0	10	10
2-Nitroaniline	U	25	1.0	25	25
Dimethyl Phthalate	U	10	1.0	10	10
2,6-Dinitrotoluene	U	10	1.0	10	10
Acenaphthylene	U	10	1.0	10	10
3-Nitroaniline	U	25	1.0	25	25
Acenaphthene	U	10	1.0	10	10
2,4-Dinitrophenol	U	25	1.0	25	25
Dibenzofuran	U	10	1.0	10	10
4-Nitrophenol	U	25	1.0	25	25
2,4-Dinitrotoluene	U	10	1.0	10	10
Diethylphthalate	U	10	1.0	10	10
Fluorene	U	10	1.0	10	10
4-Chlorophenyl-phenylether	U	10	1.0	10	10
4-Nitroaniline	U	25	1.0	25	25
4,6-Dinitro-2-Methylphenol	U	25	1.0	25	25
N-Nitrosodiphenylamine	U	10	1.0	10	10

KATAHDIN ANALYTICAL SERVICES
Report of Analytical Results

Client: Woodard & Curran
Project: Old Town
PO No:
Sample Date: 05/30/06
Received Date: 05/31/06
Extraction Date: 05/31/06
Analysis Date: 02-JUN-2006 08:15
Report Date: 06/06/2006
Matrix: WATER
% Solids: NA

Lab ID: WW2637-4
Client ID: SB-26
SDG: WW2637
Extracted by: GN
Extraction Method: SW846 3510
Analyst: LRS
Analysis Method: SW846 8270C
Lab Prep Batch: WG28661
Units: ug/L

Compound	Flags	Results	DF	PQL	Adj.PQL
4-Bromophenyl-phenylether	U	10	1.0	10	10
Hexachlorobenzene	U	10	1.0	10	10
Pentachlorophenol	U	25	1.0	25	25
Phenanthrene	U	10	1.0	10	10
Anthracene	U	10	1.0	10	10
Carbazole	U	10	1.0	10	10
Di-n-butylphthalate	U	10	1.0	10	10
Fluoranthene	U	10	1.0	10	10
Pyrene	U	10	1.0	10	10
Butylbenzylphthalate	U	10	1.0	10	10
Benzo(a)anthracene	U	10	1.0	10	10
3,3'-Dichlorobenzidine	U	10	1.0	10	10
Chrysene	U	10	1.0	10	10
bis(2-Ethylhexyl)phthalate	U	10	1.0	10	10
Di-n-octylphthalate	U	10	1.0	10	10
Benzo(b)fluoranthene	U	10	1.0	10	10
Benzo(k)fluoranthene	U	10	1.0	10	10
Benzo(a)pyrene	U	10	1.0	10	10
Indeno(1,2,3-cd)pyrene	U	10	1.0	10	10
Dibenzo(a,h)anthracene	U	10	1.0	10	10
Benzo(g,h,i)perylene	U	10	1.0	10	10
2-Fluorophenol		43%			
Phenol-D6		38%			
Nitrobenzene-D5		54%			
2-Fluorobiphenyl		62%			
2,4,6-Tribromophenol		72%			
Terphenyl-D14		89%			

Katahdin
ANALYTICAL SERVICES

REPORT OF ANALYTICAL RESULTS

Client: Erika Lloyd
Woodard & Curran
41 Hutchins Drive
Portland, ME 04102

Lab Sample ID: WW2637-004
Report Date: 6/7/2006
PO No.:
Project: Old Town

Sample Description	Matrix	Filtered	Date Sampled	Date Received
SB-26	AQ	No(Total)	05/30/2006	05/31/2006

Parameter	Result	Units	Adjusted PQL	Dilution Factor	PQL	Analytical Method	Analysis Date	By	Prep Method	Prepped Date	By	QC	Notes
ARSENIC	U 0.0050	mg/L	0.0050	5	0.001	SW846 6020	6/2/06	EAM	3010	5/31/06	MW	WE31ICW1	1
BARIUM	0.0284	mg/L	0.0050	1	0.005	SW846 6010	5/31/06	SJC	SW846 3010	5/31/06	MW	WE31ICW1	
CADMIUM	U 0.0100	mg/L	0.0100	1	0.01	SW846 6010	5/31/06	SJC	SW846 3010	5/31/06	MW	WE31ICW1	
CHROMIUM	0.0035	mg/L	0.0030	5	0.0006	SW846 6020	6/2/06	EAM	3010	5/31/06	MW	WE31ICW1	
LEAD	U 0.0010	mg/L	0.0010	5	0.0002	SW846 6020	6/2/06	EAM	3010	5/31/06	MW	WE31ICW1	1
MERCURY	U 0.20	ug/L	0.20	1	0.2	SW846 7470	6/1/06	MW	SW846 7470	6/1/06	MW	WF01HGW0	
SELENIUM	U 0.010	mg/L	0.010	1	0.01	SW846 6010	5/31/06	SJC	SW846 3010	5/31/06	MW	WE31ICW1	
SILVER	U 0.015	mg/L	0.015	1	0.015	SW846 6010	5/31/06	SJC	SW846 3010	5/31/06	MW	WE31ICW1	

1 The laboratory's Practical Quantitation Level could not be achieved for this parameter due to sample composition, matrix effects,sample volume, or quantity used for analysis.

KATAHDIN ANALYTICAL SERVICES
Report of Analytical Results

Client: Woodard & Curran
Project: Old Town
PO No:
Sample Date: 05/30/06
Received Date: 05/31/06
Extraction Date:
Analysis Date: 31-MAY-2006 17:40
Report Date: 06/02/2006
Matrix: WATER
% Solids: NA

Lab ID: WW2637-5
Client ID: SB-100
SDG: WW2637
Extracted by:
Extraction Method: SW846 5030
Analyst: SKT
Analysis Method: SW846 8260B
Lab Prep Batch: WG28646
Units: ug/l

Compound	Flags	Results	DF	PQL	Adj.PQL
Dichlorodifluoromethane	U	10	1.0	10	10
Chloromethane	U	10	1.0	10	10
Vinyl chloride	U	10	1.0	10	10
Bromomethane	U	10	1.0	10	10
Chloroethane	U	10	1.0	10	10
Trichlorofluoromethane	U	10	1.0	10	10
1,1-Dichloroethene	U	5	1.0	5	5
Methylene Chloride	U	5	1.0	5	5
trans-1,2-Dichloroethene	U	5	1.0	5	5
1,1-Dichloroethane	U	5	1.0	5	5
cis-1,2-Dichloroethene	U	5	1.0	5	5
1,2-Dichloroethylene (total)	U	10	1.0	10	10
2,2-Dichloropropane	U	5	1.0	5	5
Chloroform	U	5	1.0	5	5
Bromochloromethane	U	5	1.0	5	5
1,1,1-Trichloroethane	U	5	1.0	5	5
1,2-Dichloroethane	U	5	1.0	5	5
1,1-Dichloropropene	U	5	1.0	5	5
Carbon Tetrachloride	U	5	1.0	5	5
Benzene	U	5	1.0	5	5
1,2-Dichloropropane	U	5	1.0	5	5
Trichloroethene	U	5	1.0	5	5
Dibromomethane	U	5	1.0	5	5
Bromodichloromethane	U	5	1.0	5	5
cis-1,3-dichloropropene	U	5	1.0	5	5
Toluene	U	5	1.0	5	5
trans-1,3-Dichloropropene	U	5	1.0	5	5
1,1,2-Trichloroethane	U	5	1.0	5	5
1,3-Dichloropropane	U	5	1.0	5	5
Dibromochloromethane	U	5	1.0	5	5
Tetrachloroethene	U	5	1.0	5	5
1,2-Dibromoethane	U	5	1.0	5	5
Chlorobenzene	U	5	1.0	5	5
1,1,1,2-Tetrachloroethane	U	5	1.0	5	5
Ethylbenzene	U	5	1.0	5	5
Bromoform	U	5	1.0	5	5
Styrene	U	5	1.0	5	5
1,1,2,2-Tetrachloroethane	U	5	1.0	5	5
1,2,3-Trichloropropane	U	5	1.0	5	5
Isopropylbenzene	U	5	1.0	5	5
Bromobenzene	U	5	1.0	5	5
2-Chlorotoluene	U	5	1.0	5	5
N-Propylbenzene	U	5	1.0	5	5

KATAHDIN ANALYTICAL SERVICES
Report of Analytical Results

Client: Woodard & Curran
Project: Old Town
PO No:
Sample Date: 05/30/06
Received Date: 05/31/06
Extraction Date:
Analysis Date: 31-MAY-2006 17:40
Report Date: 06/02/2006
Matrix: WATER
% Solids: NA

Lab ID: WW2637-5
Client ID: SB-100
SDG: WW2637
Extracted by:
Extraction Method: SW846 5030
Analyst: SKT
Analysis Method: SW846 8260B
Lab Prep Batch: WG28646
Units: ug/l

Compound	Flags	Results	DF	PQL	Adj.PQL
4-Chlorotoluene	U	5	1.0	5	5
1,3,5-Trimethylbenzene	U	5	1.0	5	5
tert-Butylbenzene	U	5	1.0	5	5
1,2,4-Trichlorobenzene	U	5	1.0	5	5
sec-Butylbenzene	U	5	1.0	5	5
1,3-Dichlorobenzene	U	5	1.0	5	5
P-Isopropyltoluene	U	5	1.0	5	5
1,4-Dichlorobenzene	U	5	1.0	5	5
1,2-Dichlorobenzene	U	5	1.0	5	5
N-Butylbenzene	U	5	1.0	5	5
1,2-Dibromo-3-Chloropropane	U	5	1.0	5	5
1,2,4-Trimethylbenzene	U	5	1.0	5	5
Naphthalene	U	5	1.0	5	5
Hexachlorobutadiene	U	5	1.0	5	5
1,2,3-Trichlorobenzene	U	5	1.0	5	5
Methyl tert-butyl ether	U	10	1.0	10	10
Acetone	U	25	1.0	25	25
2-Butanone	U	25	1.0	25	25
4-methyl-2-pentanone	U	25	1.0	25	25
2-Hexanone	U	25	1.0	25	25
m+p-Xylenes	U	10	1.0	10	10
o-Xylene	U	5	1.0	5	5
Xylenes (total)	U	15	1.0	15	15
1,3,5-Trichlorobenzene	U	5	1.0	5	5
Vinyl Acetate	U	5	1.0	5	5
Carbon Disulfide	U	5	1.0	5	5
Diethyl Ether	U	5	1.0	5	5
Tetrahydrofuran	U	25	1.0	25	25
Dibromofluoromethane		92%			
1,2-Dichloroethane-D4		106%			
Toluene-D8		87%			
P-Bromofluorobenzene		88%			

KATAHDIN ANALYTICAL SERVICES
Report of Analytical Results

Client: Woodard & Curran
Project: Old Town
PO No:
Sample Date: 05/30/06
Received Date: 05/31/06
Extraction Date: 05/31/06
Analysis Date: 02-JUN-2006 08:59
Report Date: 06/06/2006
Matrix: WATER
% Solids: NA

Lab ID: WW2637-5
Client ID: SB-100
SDG: WW2637
Extracted by: GN
Extraction Method: SW846 3510
Analyst: LRS
Analysis Method: SW846 8270C
Lab Prep Batch: WG28661
Units: ug/L

Compound	Flags	Results	DF	PQL	Adj.PQL
Phenol	U	10	1.0	10	10
Bis(2-Chloroethyl)ether	U	10	1.0	10	10
2-Chlorophenol	U	10	1.0	10	10
1,3-Dichlorobenzene	U	10	1.0	10	10
1,4-Dichlorobenzene	U	10	1.0	10	10
1,2-Dichlorobenzene	U	10	1.0	10	10
2-Methylphenol	U	10	1.0	10	10
2,2'-Oxybis(1-Chloropropane)	U	10	1.0	10	10
N-Nitroso-di-n-propylamine	U	10	1.0	10	10
3&4-Methylphenol	U	10	1.0	10	10
Hexachloroethane	U	10	1.0	10	10
Nitrobenzene	U	10	1.0	10	10
Isophorone	U	10	1.0	10	10
2-Nitrophenol	U	10	1.0	10	10
2,4-Dimethylphenol	U	10	1.0	10	10
Bis(2-Chloroethoxy)methane	U	10	1.0	10	10
2,4-Dichlorophenol	U	10	1.0	10	10
1,2,4-Trichlorobenzene	U	10	1.0	10	10
Naphthalene	U	10	1.0	10	10
4-Chloroaniline	U	10	1.0	10	10
Hexachlorobutadiene	U	10	1.0	10	10
4-Chloro-3-Methylphenol	U	10	1.0	10	10
2-Methylnaphthalene	U	10	1.0	10	10
Hexachlorocyclopentadiene	U	10	1.0	10	10
2,4,6-Trichlorophenol	U	10	1.0	10	10
2,4,5-Trichlorophenol	U	25	1.0	25	25
2-Chloronaphthalene	U	10	1.0	10	10
2-Nitroaniline	U	25	1.0	25	25
Dimethyl Phthalate	U	10	1.0	10	10
2,6-Dinitrotoluene	U	10	1.0	10	10
Acenaphthylene	U	10	1.0	10	10
3-Nitroaniline	U	25	1.0	25	25
Acenaphthene	U	10	1.0	10	10
2,4-Dinitrophenol	U	25	1.0	25	25
Dibenzofuran	U	10	1.0	10	10
4-Nitrophenol	U	25	1.0	25	25
2,4-Dinitrotoluene	U	10	1.0	10	10
Diethylphthalate	U	10	1.0	10	10
Fluorene	U	10	1.0	10	10
4-Chlorophenyl-phenylether	U	10	1.0	10	10
4-Nitroaniline	U	25	1.0	25	25
4,6-Dinitro-2-Methylphenol	U	25	1.0	25	25
N-Nitrosodiphenylamine	U	10	1.0	10	10

KATAHDIN ANALYTICAL SERVICES
Report of Analytical Results

Client: Woodard & Curran
Project: Old Town
PO No:
Sample Date: 05/30/06
Received Date: 05/31/06
Extraction Date: 05/31/06
Analysis Date: 02-JUN-2006 08:59
Report Date: 06/06/2006
Matrix: WATER
% Solids: NA

Lab ID: WW2637-5
Client ID: SB-100
SDG: WW2637
Extracted by: GN
Extraction Method: SW846 3510
Analyst: LRS
Analysis Method: SW846 8270C
Lab Prep Batch: WG28661
Units: ug/L

Compound	Flags	Results	DF	PQL	Adj.PQL
4-Bromophenyl-phenylether	U	10	1.0	10	10
Hexachlorobenzene	U	10	1.0	10	10
Pentachlorophenol	U	25	1.0	25	25
Phenanthrene	U	10	1.0	10	10
Anthracene	U	10	1.0	10	10
Carbazole	U	10	1.0	10	10
Di-n-butylphthalate	U	10	1.0	10	10
Fluoranthene	U	10	1.0	10	10
Pyrene	U	10	1.0	10	10
Butylbenzylphthalate	U	10	1.0	10	10
Benzo(a)anthracene	U	10	1.0	10	10
3,3'-Dichlorobenzidine	U	10	1.0	10	10
Chrysene	U	10	1.0	10	10
bis(2-Ethylhexyl)phthalate	U	10	1.0	10	10
Di-n-octylphthalate	U	10	1.0	10	10
Benzo(b)fluoranthene	U	10	1.0	10	10
Benzo(k)fluoranthene	U	10	1.0	10	10
Benzo(a)pyrene	U	10	1.0	10	10
Indeno(1,2,3-cd)pyrene	U	10	1.0	10	10
Dibenzo(a,h)anthracene	U	10	1.0	10	10
Benzo(g,h,i)perylene	U	10	1.0	10	10
2-Fluorophenol		52%			
Phenol-D6		48%			
Nitrobenzene-D5		69%			
2-Fluorobiphenyl		74%			
2,4,6-Tribromophenol		72%			
Terphenyl-D14		95%			

Katahdin
ANALYTICAL SERVICES

REPORT OF ANALYTICAL RESULTS

Client: Erika Lloyd
Woodard & Curran
41 Hutchins Drive
Portland, ME 04102

Lab Sample ID: WW2637-005
Report Date: 6/7/2006
PO No.:
Project: Old Town

Sample Description	Matrix	Filtered	Date Sampled	Date Received
SB-100	AQ	No(Total)	05/30/2006	05/31/2006

Parameter	Result	Units	Adjusted PQL	Dilution Factor	PQL	Analytical Method	Analysis Date	By	Prep Method	Prepped Date	By	QC	Notes
ARSENIC	U 0.0050	mg/L	0.0050	5	0.001	SW846 6020	6/2/06	EAM	SW846 3010	5/31/06	MW	WE31ICW1	1
BARIUM	0.0764	mg/L	0.0050	1	0.005	SW846 6010	5/31/06	SJC	SW846 3010	5/31/06	MW	WE31ICW1	
CADMIUM	U 0.0100	mg/L	0.0100	1	0.01	SW846 6010	5/31/06	SJC	SW846 3010	5/31/06	MW	WE31ICW1	
CHROMIUM	U 0.0030	mg/L	0.0030	5	0.0006	SW846 6020	6/2/06	EAM	SW846 3010	5/31/06	MW	WE31ICW1	1
LEAD	U 0.0010	mg/L	0.0010	5	0.0002	SW846 6020	6/2/06	EAM	SW846 3010	5/31/06	MW	WE31ICW1	1
MERCURY	U 0.20	ug/L	0.20	1	0.2	SW846 7470	6/1/06	MW	SW846 7470	6/1/06	MW	WF01HGW0	
SELENIUM	U 0.010	mg/L	0.010	1	0.01	SW846 6010	5/31/06	SJC	SW846 3010	5/31/06	MW	WE31ICW1	
SILVER	U 0.015	mg/L	0.015	1	0.015	SW846 6010	5/31/06	SJC	SW846 3010	5/31/06	MW	WE31ICW1	

1 The laboratory's Practical Quantitation Level could not be achieved for this parameter due to sample composition, matrix effects,sample volume, or quantity used for analysis.

KATAHDIN ANALYTICAL SERVICES
Report of Analytical Results

Client: Woodard & Curran
Project: Old Town
PO No:
Sample Date: 05/30/06
Received Date: 05/31/06
Extraction Date:
Analysis Date: 31-MAY-2006 11:57
Report Date: 06/02/2006
Matrix: SOIL
% Solids: 44.9

Lab ID: WW2637-6
Client ID: SS01
SDG: WW2637
Extracted by:
Extraction Method: SW846 5035
Analyst: ALH
Analysis Method: SW846 8260B
Lab Prep Batch: WG28636
Units: ug/Kg

Compound	Flags	Results	DF	PQL	Adj.PQL
Dichlorodifluoromethane	U	11	1.0	10	11
Chloromethane	U	11	1.0	10	11
Vinyl chloride	U	11	1.0	10	11
Bromomethane	U	11	1.0	10	11
Chloroethane	U	11	1.0	10	11
Trichlorofluoromethane	U	11	1.0	10	11
1,1-Dichloroethene	U	6	1.0	5	6
Methylene Chloride	U	28	1.0	25	28
trans-1,2-Dichloroethene	U	6	1.0	5	6
1,1-Dichloroethane	U	6	1.0	5	6
cis-1,2-Dichloroethene	U	6	1.0	5	6
1,2-Dichloroethylene (total)	U	11	1.0	10	11
2,2-Dichloropropane	U	6	1.0	5	6
Chloroform	U	6	1.0	5	6
Bromochloromethane	U	6	1.0	5	6
1,1,1-Trichloroethane	U	6	1.0	5	6
1,2-Dichloroethane	U	6	1.0	5	6
1,1-Dichloropropene	U	6	1.0	5	6
Carbon Tetrachloride	U	6	1.0	5	6
Benzene	U	6	1.0	5	6
1,2-Dichloropropane	U	6	1.0	5	6
Trichloroethene	U	6	1.0	5	6
Dibromomethane	U	6	1.0	5	6
Bromodichloromethane	U	6	1.0	5	6
cis-1,3-dichloropropene	U	6	1.0	5	6
Toluene	U	6	1.0	5	6
trans-1,3-Dichloropropene	U	6	1.0	5	6
1,1,2-Trichloroethane	U	6	1.0	5	6
1,3-Dichloropropane	U	6	1.0	5	6
Dibromochloromethane	U	6	1.0	5	6
Tetrachloroethene	U	6	1.0	5	6
1,2-Dibromoethane	U	6	1.0	5	6
Chlorobenzene	U	6	1.0	5	6
1,1,1,2-Tetrachloroethane	U	6	1.0	5	6
Ethylbenzene	U	6	1.0	5	6
Bromoform	U	6	1.0	5	6
Styrene	U	6	1.0	5	6
1,1,2,2-Tetrachloroethane	U	6	1.0	5	6
1,2,3-Trichloropropane	U	6	1.0	5	6
Isopropylbenzene	U	6	1.0	5	6
Bromobenzene	U	6	1.0	5	6
2-Chlorotoluene	U	6	1.0	5	6
N-Propylbenzene	U	6	1.0	5	6

KATAHDIN ANALYTICAL SERVICES
Report of Analytical Results

Client: Woodard & Curran
Project: Old Town
PO No:
Sample Date: 05/30/06
Received Date: 05/31/06
Extraction Date:
Analysis Date: 31-MAY-2006 11:57
Report Date: 06/02/2006
Matrix: SOIL
% Solids: 44.9

Lab ID: WW2637-6
Client ID: SS01
SDG: WW2637
Extracted by:
Extraction Method: SW846 5035
Analyst: ALH
Analysis Method: SW846 8260B
Lab Prep Batch: WG28636
Units: ug/Kg

Compound	Flags	Results	DF	PQL	Adj.PQL
4-Chlorotoluene	U	6	1.0	5	6
1,3,5-Trimethylbenzene	U	6	1.0	5	6
tert-Butylbenzene	U	6	1.0	5	6
1,2,4-Trichlorobenzene	U	6	1.0	5	6
sec-Butylbenzene	U	6	1.0	5	6
1,3-Dichlorobenzene	U	6	1.0	5	6
P-Isopropyltoluene	U	6	1.0	5	6
1,4-Dichlorobenzene	U	6	1.0	5	6
1,2-Dichlorobenzene	U	6	1.0	5	6
N-Butylbenzene	U	6	1.0	5	6
1,2-Dibromo-3-Chloropropane	U	6	1.0	5	6
1,2,4-Trimethylbenzene	U	6	1.0	5	6
Naphthalene	U	6	1.0	5	6
Hexachlorobutadiene	U	6	1.0	5	6
1,2,3-Trichlorobenzene	U	6	1.0	5	6
Methyl tert-butyl ether	U	11	1.0	10	11
Acetone	U	28	1.0	25	28
2-Butanone	U	28	1.0	25	28
4-methyl-2-pentanone	U	28	1.0	25	28
2-Hexanone	U	28	1.0	25	28
m+p-Xylenes	U	11	1.0	10	11
o-Xylene	U	6	1.0	5	6
Xylenes (total)	U	17	1.0	15	17
1,3,5-Trichlorobenzene	U	6	1.0	5	6
Vinyl Acetate	U	6	1.0	5	6
Carbon Disulfide	U	6	1.0	5	6
Diethyl Ether	U	6	1.0	5	6
Tetrahydrofuran	U	57	1.0	50	57
Dibromofluoromethane		66%			
1,2-Dichloroethane-D4		63%			
Toluene-D8		63%			
P-Bromofluorobenzene		38%			

KATAHDIN ANALYTICAL SERVICES
Report of Analytical Results

Client: Woodard & Curran
Project: Old Town
PO No:
Sample Date: 05/30/06
Received Date: 05/31/06
Extraction Date: 05/31/06
Analysis Date: 31-MAY-2006 17:13
Report Date: 06/06/2006
Matrix: SOIL
% Solids: 44.9

Lab ID: WW2637-6
Client ID: SS01
SDG: WW2637
Extracted by: KM
Extraction Method: SW846 3550
Analyst: LRS
Analysis Method: SW846 8270C
Lab Prep Batch: WG28621
Units: ug/Kg

Compound	Flags	Results	DF	PQL	Adj.PQL
Phenol	U	740	1.0	330	740
Bis(2-Chloroethyl)ether	U	740	1.0	330	740
2-Chlorophenol	U	740	1.0	330	740
1,3-Dichlorobenzene	U	740	1.0	330	740
1,4-Dichlorobenzene	U	740	1.0	330	740
1,2-Dichlorobenzene	U	740	1.0	330	740
2-Methylphenol	U	740	1.0	330	740
2,2'-Oxybis(1-Chloropropane)	U	740	1.0	330	740
N-Nitroso-di-n-propylamine	U	740	1.0	330	740
3&4-Methylphenol		2400	1.0	330	740
Hexachloroethane	U	740	1.0	330	740
Nitrobenzene	U	740	1.0	330	740
Isophorone	U	740	1.0	330	740
2-Nitrophenol	U	740	1.0	330	740
2,4-Dimethylphenol	U	740	1.0	330	740
Bis(2-Chloroethoxy)methane	U	740	1.0	330	740
2,4-Dichlorophenol	U	740	1.0	330	740
1,2,4-Trichlorobenzene	U	740	1.0	330	740
Naphthalene	U	740	1.0	330	740
4-Chloroaniline	U	740	1.0	330	740
Hexachlorobutadiene	U	740	1.0	330	740
4-Chloro-3-Methylphenol	U	740	1.0	330	740
2-Methylnaphthalene	U	740	1.0	330	740
Hexachlorocyclopentadiene	U	740	1.0	330	740
2,4,6-Trichlorophenol	U	740	1.0	330	740
2,4,5-Trichlorophenol	U	1800	1.0	820	1800
2-Chloronaphthalene	U	740	1.0	330	740
2-Nitroaniline	U	1800	1.0	820	1800
Dimethyl Phthalate	U	740	1.0	330	740
2,6-Dinitrotoluene	U	740	1.0	330	740
Acenaphthylene	U	740	1.0	330	740
3-Nitroaniline	U	1800	1.0	820	1800
Acenaphthene	U	740	1.0	330	740
2,4-Dinitrophenol	U	1800	1.0	820	1800
Dibenzofuran	U	740	1.0	330	740
4-Nitrophenol	U	1800	1.0	820	1800
2,4-Dinitrotoluene	U	1800	1.0	820	1800
Diethylphthalate	U	740	1.0	330	740
Fluorene	U	740	1.0	330	740
4-Chlorophenyl-phenylether	U	740	1.0	330	740
4-Nitroaniline	U	1800	1.0	820	1800
4,6-Dinitro-2-Methylphenol	U	1800	1.0	820	1800
N-Nitrosodiphenylamine	U	740	1.0	330	740

KATAHDIN ANALYTICAL SERVICES
Report of Analytical Results

Client: Woodard & Curran
Project: Old Town
PO No:
Sample Date: 05/30/06
Received Date: 05/31/06
Extraction Date: 05/31/06
Analysis Date: 31-MAY-2006 17:13
Report Date: 06/06/2006
Matrix: SOIL
% Solids: 44.9

Lab ID: WW2637-6
Client ID: SS01
SDG: WW2637
Extracted by: KM
Extraction Method: SW846 3550
Analyst: LRS
Analysis Method: SW846 8270C
Lab Prep Batch: WG28621
Units: ug/Kg

Compound	Flags	Results	DF	PQL	Adj.PQL
4-Bromophenyl-phenylether	U	740	1.0	330	740
Hexachlorobenzene	U	740	1.0	330	740
Pentachlorophenol	U	1800	1.0	820	1800
Phenanthrene	U	740	1.0	330	740
Anthracene	U	740	1.0	330	740
Carbazole	U	740	1.0	330	740
Di-n-butylphthalate	U	740	1.0	330	740
Fluoranthene	U	740	1.0	330	740
Pyrene	U	740	1.0	330	740
Butylbenzylphthalate	U	740	1.0	330	740
Benzo(a)anthracene	U	740	1.0	330	740
3,3'-Dichlorobenzidine	U	740	1.0	330	740
Chrysene	U	740	1.0	330	740
bis(2-Ethylhexyl)phthalate	U	740	1.0	330	740
Di-n-octylphthalate	U	740	1.0	330	740
Benzo(b)fluoranthene	U	740	1.0	330	740
Benzo(k)fluoranthene	U	740	1.0	330	740
Benzo(a)pyrene	U	740	1.0	330	740
Indeno(1,2,3-cd)pyrene	U	740	1.0	330	740
Dibenzo(a,h)anthracene	U	740	1.0	330	740
Benzo(g,h,i)perylene	U	740	1.0	330	740
2-Fluorophenol		53%			
Phenol-D6		60%			
Nitrobenzene-D5		52%			
2-Fluorobiphenyl		57%			
2,4,6-Tribromophenol		66%			
Terphenyl-D14		72%			

Katahdin
ANALYTICAL SERVICES

REPORT OF ANALYTICAL RESULTS

Client: Erika Lloyd
Woodard & Curran
41 Hutchins Drive
Portland, ME 04102

Lab Sample ID: WW2637-006
Report Date: 6/7/2006
PO No.:
Project: Old Town

Sample Description	Matrix	Percent Solids(%)	Date Sampled	Date Received
SS01	SL	44.9	05/30/2006	05/31/2006

Parameter	Result	Units	Adjusted PQL	Dilution Factor	PQL	Analytical Method	Analysis Date	By	Prep Method	Prepped Date	By	QC	Notes
ARSENIC	2.3	mg/Kg	2.	1	0.8	SW846 6010	6/1/06	SJC	SW846 3050	5/31/06	EAM	WE31ICS0	
BARIUM	149.	mg/Kg	1.0	1	0.5	SW846 6010	6/1/06	SJC	SW846 3050	5/31/06	EAM	WE31ICS0	
CADMIUM	0.634	mg/Kg	0.206	5	0.02	SW846 6020	6/2/06	EAM	SW846 3050	5/31/06	EAM	WE31ICS0	
CHROMIUM	11.8	mg/Kg	0.62	5	0.06	SW846 6020	6/2/06	EAM	SW846 3050	5/31/06	EAM	WE31ICS0	
LEAD	5.50	mg/Kg	0.21	5	0.02	SW846 6020	6/2/06	EAM	SW846 3050	5/31/06	EAM	WE31ICS0	
MERCURY	U 0.077	ug/g	0.077	1	0.04	SW846 7471	6/2/06	MW	SW846 7471	6/1/06	MW	WF01HGS0	1
SELENIUM	U 2.1	mg/Kg	2.1	1	1	SW846 6010	6/1/06	SJC	SW846 3050	5/31/06	EAM	WE31ICS0	1
SILVER	U 3.1	mg/Kg	3.1	1	1.5	SW846 6010	6/1/06	SJC	SW846 3050	5/31/06	EAM	WE31ICS0	1

1 The laboratory's Practical Quantitation Level could not be achieved for this parameter due to sample composition, matrix effects,sample volume, or quantity used for analysis.

KATAHDIN ANALYTICAL SERVICES
Report of Analytical Results

Client: Woodard & Curran
Project: Old Town
PO No:
Sample Date: 05/30/06
Received Date: 05/31/06
Extraction Date: 05/31/06
Analysis Date: 02-JUN-2006 10:31
Report Date: 06/06/2006
Matrix: SOIL
% Solids: 92.2

Lab ID: WW2637-7DL2
Client ID: SS02
SDG: WW2637
Extracted by: KM
Extraction Method: SW846 3550
Analyst: LRS
Analysis Method: SW846 8270C
Lab Prep Batch: WG28621
Units: ug/Kg

Compound	Flags	Results	DF	PQL	Adj.PQL
Phenol	U	1800	5.0	330	1800
Bis(2-Chloroethyl)ether	U	1800	5.0	330	1800
2-Chlorophenol	U	1800	5.0	330	1800
1,3-Dichlorobenzene	U	1800	5.0	330	1800
1,4-Dichlorobenzene	U	1800	5.0	330	1800
1,2-Dichlorobenzene	U	1800	5.0	330	1800
2-Methylphenol	U	1800	5.0	330	1800
2,2'-Oxybis(1-Chloropropane)	U	1800	5.0	330	1800
N-Nitroso-di-n-propylamine	U	1800	5.0	330	1800
3&4-Methylphenol	U	1800	5.0	330	1800
Hexachloroethane	U	1800	5.0	330	1800
Nitrobenzene	U	1800	5.0	330	1800
Isophorone	U	1800	5.0	330	1800
2-Nitrophenol	U	1800	5.0	330	1800
2,4-Dimethylphenol	U	1800	5.0	330	1800
Bis(2-Chloroethoxy)methane	U	1800	5.0	330	1800
2,4-Dichlorophenol	U	1800	5.0	330	1800
1,2,4-Trichlorobenzene	U	1800	5.0	330	1800
Naphthalene	U	1800	5.0	330	1800
4-Chloroaniline	U	1800	5.0	330	1800
Hexachlorobutadiene	U	1800	5.0	330	1800
4-Chloro-3-Methylphenol	U	1800	5.0	330	1800
2-Methylnaphthalene		6200	5.0	330	1800
Hexachlorocyclopentadiene	U	1800	5.0	330	1800
2,4,6-Trichlorophenol	U	1800	5.0	330	1800
2,4,5-Trichlorophenol	U	4400	5.0	820	4400
2-Chloronaphthalene	U	1800	5.0	330	1800
2-Nitroaniline	U	4400	5.0	820	4400
Dimethyl Phthalate	U	1800	5.0	330	1800
2,6-Dinitrotoluene	U	1800	5.0	330	1800
Acenaphthylene	U	1800	5.0	330	1800
3-Nitroaniline	U	4400	5.0	820	4400
Acenaphthene	U	1800	5.0	330	1800
2,4-Dinitrophenol	U	4400	5.0	820	4400
Dibenzofuran	U	1800	5.0	330	1800
4-Nitrophenol	U	4400	5.0	820	4400
2,4-Dinitrotoluene	U	4400	5.0	820	4400
Diethylphthalate	U	1800	5.0	330	1800
Fluorene	U	1800	5.0	330	1800
4-Chlorophenyl-phenylether	U	1800	5.0	330	1800
4-Nitroaniline	U	4400	5.0	820	4400
4,6-Dinitro-2-Methylphenol	U	4400	5.0	820	4400
N-Nitrosodiphenylamine	U	1800	5.0	330	1800

KATAHDIN ANALYTICAL SERVICES
Report of Analytical Results

Client: Woodard & Curran
Project: Old Town
PO No:
Sample Date: 05/30/06
Received Date: 05/31/06
Extraction Date: 05/31/06
Analysis Date: 02-JUN-2006 10:31
Report Date: 06/06/2006
Matrix: SOIL
% Solids: 92.2

Lab ID: WW2637-7DL2
Client ID: SS02
SDG: WW2637
Extracted by: KM
Extraction Method: SW846 3550
Analyst: LRS
Analysis Method: SW846 8270C
Lab Prep Batch: WG28621
Units: ug/Kg

Compound	Flags	Results	DF	PQL	Adj.PQL
4-Bromophenyl-phenylether	U	1800	5.0	330	1800
Hexachlorobenzene	U	1800	5.0	330	1800
Pentachlorophenol	U	4400	5.0	820	4400
Phenanthrene		3300	5.0	330	1800
Anthracene	U	1800	5.0	330	1800
Carbazole	U	1800	5.0	330	1800
Di-n-butylphthalate	U	1800	5.0	330	1800
Fluoranthene	U	1800	5.0	330	1800
Pyrene	U	1800	5.0	330	1800
Butylbenzylphthalate	U	1800	5.0	330	1800
Benzo(a)anthracene	U	1800	5.0	330	1800
3,3'-Dichlorobenzidine	U	1800	5.0	330	1800
Chrysene	U	1800	5.0	330	1800
bis(2-Ethylhexyl)phthalate	U	1800	5.0	330	1800
Di-n-octylphthalate	U	1800	5.0	330	1800
Benzo(b)fluoranthene	U	1800	5.0	330	1800
Benzo(k)fluoranthene	U	1800	5.0	330	1800
Benzo(a)pyrene	U	1800	5.0	330	1800
Indeno(1,2,3-cd)pyrene	U	1800	5.0	330	1800
Dibenzo(a,h)anthracene	U	1800	5.0	330	1800
Benzo(g,h,i)perylene	U	1800	5.0	330	1800
2-Fluorophenol		56%			
Phenol-D6		65%			
Nitrobenzene-D5		89%			
2-Fluorobiphenyl		82%			
2,4,6-Tribromophenol		65%			
Terphenyl-D14		117%			

KATAHDIN ANALYTICAL SERVICES
Report of Analytical Results

Client: Woodard & Curran
Project: Old Town
PO No:
Sample Date: 05/30/06
Received Date: 05/31/06
Extraction Date:
Analysis Date: 31-MAY-2006 15:01
Report Date: 06/02/2006
Matrix: SOIL
% Solids: 53.3

Lab ID: WW2637-8RA
Client ID: SS03
SDG: WW2637
Extracted by:
Extraction Method: SW846 5035
Analyst: ALH
Analysis Method: SW846 8260B
Lab Prep Batch: WG28636
Units: ug/Kg

Compound	Flags	Results	DF	PQL	Adj.PQL
Dichlorodifluoromethane	U	11	1.0	10	11
Chloromethane	U	11	1.0	10	11
Vinyl chloride	U	11	1.0	10	11
Bromomethane	U	11	1.0	10	11
Chloroethane	U	11	1.0	10	11
Trichlorofluoromethane	U	11	1.0	10	11
1,1-Dichloroethene	U	5	1.0	5	5
Methylene Chloride	U	27	1.0	25	27
trans-1,2-Dichloroethene	U	5	1.0	5	5
1,1-Dichloroethane	U	5	1.0	5	5
cis-1,2-Dichloroethene	U	5	1.0	5	5
1,2-Dichloroethylene (total)	U	11	1.0	10	11
2,2-Dichloropropane	U	5	1.0	5	5
Chloroform	U	5	1.0	5	5
Bromochloromethane	U	5	1.0	5	5
1,1,1-Trichloroethane	U	5	1.0	5	5
1,2-Dichloroethane	U	5	1.0	5	5
1,1-Dichloropropene	U	5	1.0	5	5
Carbon Tetrachloride	U	5	1.0	5	5
Benzene	U	5	1.0	5	5
1,2-Dichloropropane	U	5	1.0	5	5
Trichloroethene	U	5	1.0	5	5
Dibromomethane	U	5	1.0	5	5
Bromodichloromethane	U	5	1.0	5	5
cis-1,3-dichloropropene	U	5	1.0	5	5
Toluene	U	5	1.0	5	5
trans-1,3-Dichloropropene	U	5	1.0	5	5
1,1,2-Trichloroethane	U	5	1.0	5	5
1,3-Dichloropropane	U	5	1.0	5	5
Dibromochloromethane	U	5	1.0	5	5
Tetrachloroethene	U	5	1.0	5	5
1,2-Dibromoethane	U	5	1.0	5	5
Chlorobenzene	U	5	1.0	5	5
1,1,1,2-Tetrachloroethane	U	5	1.0	5	5
Ethylbenzene	U	5	1.0	5	5
Bromoform	U	5	1.0	5	5
Styrene	U	5	1.0	5	5
1,1,2,2-Tetrachloroethane	U	5	1.0	5	5
1,2,3-Trichloropropane	U	5	1.0	5	5
Isopropylbenzene	U	5	1.0	5	5
Bromobenzene	U	5	1.0	5	5
2-Chlorotoluene	U	5	1.0	5	5
N-Propylbenzene	U	5	1.0	5	5

KATAHDIN ANALYTICAL SERVICES
Report of Analytical Results

Client: Woodard & Curran
Project: Old Town
PO No:
Sample Date: 05/30/06
Received Date: 05/31/06
Extraction Date:
Analysis Date: 31-MAY-2006 15:01
Report Date: 06/02/2006
Matrix: SOIL
% Solids: 53.3

Lab ID: WW2637-8RA
Client ID: SS03
SDG: WW2637
Extracted by:
Extraction Method: SW846 5035
Analyst: ALH
Analysis Method: SW846 8260B
Lab Prep Batch: WG28636
Units: ug/Kg

Compound	Flags	Results	DF	PQL	Adj.PQL
4-Chlorotoluene	U	5	1.0	5	5
1,3,5-Trimethylbenzene	U	5	1.0	5	5
tert-Butylbenzene	U	5	1.0	5	5
1,2,4-Trichlorobenzene	U	5	1.0	5	5
sec-Butylbenzene	U	5	1.0	5	5
1,3-Dichlorobenzene	U	5	1.0	5	5
P-Isopropyltoluene	U	5	1.0	5	5
1,4-Dichlorobenzene	U	5	1.0	5	5
1,2-Dichlorobenzene	U	5	1.0	5	5
N-Butylbenzene	U	5	1.0	5	5
1,2-Dibromo-3-Chloropropane	U	5	1.0	5	5
1,2,4-Trimethylbenzene	U	5	1.0	5	5
Naphthalene	U	5	1.0	5	5
Hexachlorobutadiene	U	5	1.0	5	5
1,2,3-Trichlorobenzene	U	5	1.0	5	5
Methyl tert-butyl ether	U	11	1.0	10	11
Acetone	U	27	1.0	25	27
2-Butanone	U	27	1.0	25	27
4-methyl-2-pentanone	U	27	1.0	25	27
2-Hexanone	U	27	1.0	25	27
m+p-Xylenes	U	11	1.0	10	11
o-Xylene	U	5	1.0	5	5
Xylenes (total)	U	16	1.0	15	16
1,3,5-Trichlorobenzene	U	5	1.0	5	5
Vinyl Acetate	U	5	1.0	5	5
Carbon Disulfide	U	5	1.0	5	5
Diethyl Ether	U	5	1.0	5	5
Tetrahydrofuran	U	54	1.0	50	54
Dibromofluoromethane		68%			
1,2-Dichloroethane-D4		70%			
Toluene-D8		54%			
P-Bromofluorobenzene		* 26%			

KATAHDIN ANALYTICAL SERVICES
Report of Analytical Results

Client: Woodard & Curran
Project: Old Town
PO No:
Sample Date: 05/30/06
Received Date: 05/31/06
Extraction Date: 05/31/06
Analysis Date: 31-MAY-2006 17:59
Report Date: 06/06/2006
Matrix: SOIL
% Solids: 53.3

Lab ID: WW2637-8
Client ID: SS03
SDG: WW2637
Extracted by: KM
Extraction Method: SW846 3550
Analyst: LRS
Analysis Method: SW846 8270C
Lab Prep Batch: WG28621
Units: ug/Kg

Compound	Flags	Results	DF	PQL	Adj.PQL
Phenol	U	620	1.0	330	620
Bis(2-Chloroethyl)ether	U	620	1.0	330	620
2-Chlorophenol	U	620	1.0	330	620
1,3-Dichlorobenzene	U	620	1.0	330	620
1,4-Dichlorobenzene	U	620	1.0	330	620
1,2-Dichlorobenzene	U	620	1.0	330	620
2-Methylphenol	U	620	1.0	330	620
2,2'-Oxybis(1-Chloropropane)	U	620	1.0	330	620
N-Nitroso-di-n-propylamine	U	620	1.0	330	620
3&4-Methylphenol	U	620	1.0	330	620
Hexachloroethane	U	620	1.0	330	620
Nitrobenzene	U	620	1.0	330	620
Isophorone	U	620	1.0	330	620
2-Nitrophenol	U	620	1.0	330	620
2,4-Dimethylphenol	U	620	1.0	330	620
Bis(2-Chloroethoxy)methane	U	620	1.0	330	620
2,4-Dichlorophenol	U	620	1.0	330	620
1,2,4-Trichlorobenzene	U	620	1.0	330	620
Naphthalene	U	620	1.0	330	620
4-Chloroaniline	U	620	1.0	330	620
Hexachlorobutadiene	U	620	1.0	330	620
4-Chloro-3-Methylphenol	U	620	1.0	330	620
2-Methylnaphthalene	U	620	1.0	330	620
Hexachlorocyclopentadiene	U	620	1.0	330	620
2,4,6-Trichlorophenol	U	620	1.0	330	620
2,4,5-Trichlorophenol	U	1500	1.0	820	1500
2-Chloronaphthalene	U	620	1.0	330	620
2-Nitroaniline	U	1500	1.0	820	1500
Dimethyl Phthalate	U	620	1.0	330	620
2,6-Dinitrotoluene	U	620	1.0	330	620
Acenaphthylene	U	620	1.0	330	620
3-Nitroaniline	U	1500	1.0	820	1500
Acenaphthene	U	620	1.0	330	620
2,4-Dinitrophenol	U	1500	1.0	820	1500
Dibenzofuran	U	620	1.0	330	620
4-Nitrophenol	U	1500	1.0	820	1500
2,4-Dinitrotoluene	U	1500	1.0	820	1500
Diethylphthalate	U	620	1.0	330	620
Fluorene	U	620	1.0	330	620
4-Chlorophenyl-phenylether	U	620	1.0	330	620
4-Nitroaniline	U	1500	1.0	820	1500
4,6-Dinitro-2-Methylphenol	U	1500	1.0	820	1500
N-Nitrosodiphenylamine	U	620	1.0	330	620

KATAHDIN ANALYTICAL SERVICES
Report of Analytical Results

Client: Woodard & Curran
Project: Old Town
PO No:
Sample Date: 05/30/06
Received Date: 05/31/06
Extraction Date: 05/31/06
Analysis Date: 31-MAY-2006 17:59
Report Date: 06/06/2006
Matrix: SOIL
% Solids: 53.3

Lab ID: WW2637-8
Client ID: SS03
SDG: WW2637
Extracted by: KM
Extraction Method: SW846 3550
Analyst: LRS
Analysis Method: SW846 8270C
Lab Prep Batch: WG28621
Units: ug/Kg

Compound	Flags	Results	DF	PQL	Adj.PQL
4-Bromophenyl-phenylether	U	620	1.0	330	620
Hexachlorobenzene	U	620	1.0	330	620
Pentachlorophenol	U	1500	1.0	820	1500
Phenanthrene	U	620	1.0	330	620
Anthracene	U	620	1.0	330	620
Carbazole	U	620	1.0	330	620
Di-n-butylphthalate	U	620	1.0	330	620
Fluoranthene	U	620	1.0	330	620
Pyrene	U	620	1.0	330	620
Butylbenzylphthalate	U	620	1.0	330	620
Benzo(a)anthracene	U	620	1.0	330	620
3,3'-Dichlorobenzidine	U	620	1.0	330	620
Chrysene	U	620	1.0	330	620
bis(2-Ethylhexyl)phthalate	U	620	1.0	330	620
Di-n-octylphthalate	U	620	1.0	330	620
Benzo(b)fluoranthene	U	620	1.0	330	620
Benzo(k)fluoranthene	U	620	1.0	330	620
Benzo(a)pyrene	U	620	1.0	330	620
Indeno(1,2,3-cd)pyrene	U	620	1.0	330	620
Dibenzo(a,h)anthracene	U	620	1.0	330	620
Benzo(g,h,i)perylene	U	620	1.0	330	620
2-Fluorophenol		52%			
Phenol-D6		61%			
Nitrobenzene-D5		53%			
2-Fluorobiphenyl		59%			
2,4,6-Tribromophenol		63%			
Terphenyl-D14		85%			

Katahdin
ANALYTICAL SERVICES

REPORT OF ANALYTICAL RESULTS

Client: Erika Lloyd
Woodard & Curran
41 Hutchins Drive
Portland, ME 04102

Lab Sample ID: WW2637-008
Report Date: 6/7/2006
PO No.:
Project: Old Town

Sample Description	Matrix	Percent Solids(%)	Date Sampled	Date Received
SS03	SL	53.3	05/30/2006	05/31/2006

Parameter	Result	Units	Adjusted PQL	Dilution Factor	PQL	Analytical Method	Analysis Date	By	Prep Method	Prepped Date	By	QC	Notes
ARSENIC	1.	mg/Kg	1.	1	0.8	SW846 6010	6/1/06	SJC	SW846 3050	5/31/06	EAM	WE31ICS0	
BARIUM	125.	mg/Kg	0.82	1	0.5	SW846 6010	6/1/06	SJC	SW846 3050	5/31/06	EAM	WE31ICS0	
CADMIUM	0.811	mg/Kg	0.164	5	0.02	SW846 6020	6/2/06	EAM	SW846 3050	5/31/06	EAM	WE31ICS0	
CHROMIUM	9.05	mg/Kg	0.49	5	0.06	SW846 6020	6/2/06	EAM	SW846 3050	5/31/06	EAM	WE31ICS0	
LEAD	8.00	mg/Kg	0.16	5	0.02	SW846 6020	6/2/06	EAM	SW846 3050	5/31/06	EAM	WE31ICS0	
MERCURY	U 0.066	ug/g	0.066	1	0.04	SW846 7471	6/2/06	MW	SW846 7471	6/1/06	MW	WF01HGS0	1
SELENIUM	U 1.6	mg/Kg	1.6	1	1	SW846 6010	6/1/06	SJC	SW846 3050	5/31/06	EAM	WE31ICS0	1
SILVER	U 2.5	mg/Kg	2.5	1	1.5	SW846 6010	6/1/06	SJC	SW846 3050	5/31/06	EAM	WE31ICS0	1

1 The laboratory's Practical Quantitation Level could not be achieved for this parameter due to sample composition, matrix effects,sample volume, or quantity used for analysis.

KATAHDIN ANALYTICAL SERVICES
Report of Analytical Results

Client: Woodard & Curran
Project: Old Town
PO No:
Sample Date: 05/30/06
Received Date: 05/31/06
Extraction Date: 05/31/06
Analysis Date: 31-MAY-2006 18:44
Report Date: 06/06/2006
Matrix: SOIL
% Solids: 50.3

Lab ID: WW2637-9
Client ID: SS80
SDG: WW2637
Extracted by: KM
Extraction Method: SW846 3550
Analyst: LRS
Analysis Method: SW846 8270C
Lab Prep Batch: WG28621
Units: ug/Kg

Compound	Flags	Results	DF	PQL	Adj.PQL
Phenol	U	660	1.0	330	660
Bis(2-Chloroethyl)ether	U	660	1.0	330	660
2-Chlorophenol	U	660	1.0	330	660
1,3-Dichlorobenzene	U	660	1.0	330	660
1,4-Dichlorobenzene	U	660	1.0	330	660
1,2-Dichlorobenzene	U	660	1.0	330	660
2-Methylphenol	U	660	1.0	330	660
2,2'-Oxybis(1-Chloropropane)	U	660	1.0	330	660
N-Nitroso-di-n-propylamine	U	660	1.0	330	660
3&4-Methylphenol	U	660	1.0	330	660
Hexachloroethane	U	660	1.0	330	660
Nitrobenzene	U	660	1.0	330	660
Isophorone	U	660	1.0	330	660
2-Nitrophenol	U	660	1.0	330	660
2,4-Dimethylphenol	U	660	1.0	330	660
Bis(2-Chloroethoxy)methane	U	660	1.0	330	660
2,4-Dichlorophenol	U	660	1.0	330	660
1,2,4-Trichlorobenzene	U	660	1.0	330	660
Naphthalene	U	660	1.0	330	660
4-Chloroaniline	U	660	1.0	330	660
Hexachlorobutadiene	U	660	1.0	330	660
4-Chloro-3-Methylphenol	U	660	1.0	330	660
2-Methylnaphthalene	U	660	1.0	330	660
Hexachlorocyclopentadiene	U	660	1.0	330	660
2,4,6-Trichlorophenol	U	660	1.0	330	660
2,4,5-Trichlorophenol	U	1600	1.0	820	1600
2-Chloronaphthalene	U	660	1.0	330	660
2-Nitroaniline	U	1600	1.0	820	1600
Dimethyl Phthalate	U	660	1.0	330	660
2,6-Dinitrotoluene	U	660	1.0	330	660
Acenaphthylene	U	660	1.0	330	660
3-Nitroaniline	U	1600	1.0	820	1600
Acenaphthene	U	660	1.0	330	660
2,4-Dinitrophenol	U	1600	1.0	820	1600
Dibenzofuran	U	660	1.0	330	660
4-Nitrophenol	U	1600	1.0	820	1600
2,4-Dinitrotoluene	U	1600	1.0	820	1600
Diethylphthalate	U	660	1.0	330	660
Fluorene	U	660	1.0	330	660
4-Chlorophenyl-phenylether	U	660	1.0	330	660
4-Nitroaniline	U	1600	1.0	820	1600
4,6-Dinitro-2-Methylphenol	U	1600	1.0	820	1600
N-Nitrosodiphenylamine	U	660	1.0	330	660

KATAHDIN ANALYTICAL SERVICES
Report of Analytical Results

Client: Woodard & Curran
Project: Old Town
PO No:
Sample Date: 05/30/06
Received Date: 05/31/06
Extraction Date: 05/31/06
Analysis Date: 31-MAY-2006 18:44
Report Date: 06/06/2006
Matrix: SOIL
% Solids: 50.3

Lab ID: WW2637-9
Client ID: SS80
SDG: WW2637
Extracted by: KM
Extraction Method: SW846 3550
Analyst: LRS
Analysis Method: SW846 8270C
Lab Prep Batch: WG28621
Units: ug/Kg

Compound	Flags	Results	DF	PQL	Adj.PQL
4-Bromophenyl-phenylether	U	660	1.0	330	660
Hexachlorobenzene	U	660	1.0	330	660
Pentachlorophenol	U	1600	1.0	820	1600
Phenanthrene	U	660	1.0	330	660
Anthracene	U	660	1.0	330	660
Carbazole	U	660	1.0	330	660
Di-n-butylphthalate	U	660	1.0	330	660
Fluoranthene	U	660	1.0	330	660
Pyrene	U	660	1.0	330	660
Butylbenzylphthalate	U	660	1.0	330	660
Benzo(a)anthracene	U	660	1.0	330	660
3,3'-Dichlorobenzidine	U	660	1.0	330	660
Chrysene	U	660	1.0	330	660
bis(2-Ethylhexyl)phthalate	U	660	1.0	330	660
Di-n-octylphthalate	U	660	1.0	330	660
Benzo(b)fluoranthene	U	660	1.0	330	660
Benzo(k)fluoranthene	U	660	1.0	330	660
Benzo(a)pyrene	U	660	1.0	330	660
Indeno(1,2,3-cd)pyrene	U	660	1.0	330	660
Dibenzo(a,h)anthracene	U	660	1.0	330	660
Benzo(g,h,i)perylene	U	660	1.0	330	660
2-Fluorophenol		53%			
Phenol-D6		64%			
Nitrobenzene-D5		57%			
2-Fluorobiphenyl		61%			
2,4,6-Tribromophenol		67%			
Terphenyl-D14		85%			

Katahdin
ANALYTICAL SERVICES

REPORT OF ANALYTICAL RESULTS

Client: Erika Lloyd
Woodard & Curran
41 Hutchins Drive
Portland, ME 04102

Lab Sample ID: WW2637-009
Report Date: 6/7/2006
PO No.:
Project: Old Town

Sample Description	Matrix	Percent Solids(%)	Date Sampled	Date Received
SS80	SL	50.3	05/30/2006	05/31/2006

Parameter	Result	Units	Adjusted PQL	Dilution Factor	PQL	Analytical Method	Analysis Date	By	Prep Method	Prepped Date	By	QC	Notes
ARSENIC	1.9	mg/Kg	1.	1	0.8	SW846 6010	6/1/06	SJC	SW846 3050	5/31/06	EAM	WE31ICS0	
BARIUM	148.	mg/Kg	0.86	1	0.5	SW846 6010	6/1/06	SJC	SW846 3050	5/31/06	EAM	WE31ICS0	
CADMIUM	0.891	mg/Kg	0.171	5	0.02	SW846 6020	6/2/06	EAM	SW846 3050	5/31/06	EAM	WE31ICS0	
CHROMIUM	9.43	mg/Kg	0.51	5	0.06	SW846 6020	6/2/06	EAM	SW846 3050	5/31/06	EAM	WE31ICS0	
LEAD	8.07	mg/Kg	0.17	5	0.02	SW846 6020	6/2/06	EAM	SW846 3050	5/31/06	EAM	WE31ICS0	
MERCURY	U 0.073	ug/g	0.073	1	0.04	SW846 7471	6/2/06	MW	SW846 7471	6/1/06	MW	WF01HGS0	1
SELENIUM	U 1.7	mg/Kg	1.7	1	1	SW846 6010	6/1/06	SJC	SW846 3050	5/31/06	EAM	WE31ICS0	1
SILVER	U 2.6	mg/Kg	2.6	1	1.5	SW846 6010	6/1/06	SJC	SW846 3050	5/31/06	EAM	WE31ICS0	1

1 The laboratory's Practical Quantitation Level could not be achieved for this parameter due to sample composition, matrix effects,sample volume, or quantity used for analysis.

KATAHDIN ANALYTICAL SERVICES
Report of Analytical Results

Client: Woodard & Curran
Project: Old Town
PO No:
Sample Date: 05/30/06
Received Date: 05/31/06
Extraction Date:
Analysis Date: 31-MAY-2006 13:10
Report Date: 06/07/2006
Matrix: SOIL
% Solids: 52.5

Lab ID: WW2637-10
Client ID: SS04
SDG: WW2637
Extracted by:
Extraction Method: SW846 5035
Analyst: ALH
Analysis Method: SW846 8260B
Lab Prep Batch: WG28636
Units: ug/Kg

Compound	Flags	Results	DF	PQL	Adj.PQL
Dichlorodifluoromethane	U	12	1.0	10	12
Chloromethane	U	12	1.0	10	12
Vinyl chloride	U	12	1.0	10	12
Bromomethane	U	12	1.0	10	12
Chloroethane	U	12	1.0	10	12
Trichlorofluoromethane	U	12	1.0	10	12
1,1-Dichloroethene	U	6	1.0	5	6
Methylene Chloride	U	31	1.0	25	31
trans-1,2-Dichloroethene	U	6	1.0	5	6
1,1-Dichloroethane	U	6	1.0	5	6
cis-1,2-Dichloroethene	U	6	1.0	5	6
1,2-Dichloroethylene (total)	U	12	1.0	10	12
2,2-Dichloropropane	U	6	1.0	5	6
Chloroform	U	6	1.0	5	6
Bromochloromethane	U	6	1.0	5	6
1,1,1-Trichloroethane	U	6	1.0	5	6
1,2-Dichloroethane	U	6	1.0	5	6
1,1-Dichloropropene	U	6	1.0	5	6
Carbon Tetrachloride	U	6	1.0	5	6
Benzene	U	6	1.0	5	6
1,2-Dichloropropane	U	6	1.0	5	6
Trichloroethene	U	6	1.0	5	6
Dibromomethane	U	6	1.0	5	6
Bromodichloromethane	U	6	1.0	5	6
cis-1,3-dichloropropene	U	6	1.0	5	6
Toluene	U	6	1.0	5	6
trans-1,3-Dichloropropene	U	6	1.0	5	6
1,1,2-Trichloroethane	U	6	1.0	5	6
1,3-Dichloropropane	U	6	1.0	5	6
Dibromochloromethane	U	6	1.0	5	6
Tetrachloroethene	U	6	1.0	5	6
1,2-Dibromoethane	U	6	1.0	5	6
Chlorobenzene	U	6	1.0	5	6
1,1,1,2-Tetrachloroethane	U	6	1.0	5	6
Ethylbenzene	U	6	1.0	5	6
Bromoform	U	6	1.0	5	6
Styrene	U	6	1.0	5	6
1,1,2,2-Tetrachloroethane	U	6	1.0	5	6
1,2,3-Trichloropropane	U	6	1.0	5	6
Isopropylbenzene	U	6	1.0	5	6
Bromobenzene	U	6	1.0	5	6
2-Chlorotoluene	U	6	1.0	5	6
N-Propylbenzene	U	6	1.0	5	6

KATAHDIN ANALYTICAL SERVICES
Report of Analytical Results

Client: Woodard & Curran
Project: Old Town
PO No:
Sample Date: 05/30/06
Received Date: 05/31/06
Extraction Date:
Analysis Date: 31-MAY-2006 13:10
Report Date: 06/07/2006
Matrix: SOIL
% Solids: 52.5

Lab ID: WW2637-10
Client ID: SS04
SDG: WW2637
Extracted by:
Extraction Method: SW846 5035
Analyst: ALH
Analysis Method: SW846 8260B
Lab Prep Batch: WG28636
Units: ug/Kg

Compound	Flags	Results	DF	PQL	Adj.PQL
4-Chlorotoluene	U	6	1.0	5	6
1,3,5-Trimethylbenzene	U	6	1.0	5	6
tert-Butylbenzene	U	6	1.0	5	6
1,2,4-Trichlorobenzene	U	6	1.0	5	6
sec-Butylbenzene	U	6	1.0	5	6
1,3-Dichlorobenzene	U	6	1.0	5	6
P-Isopropyltoluene	U	6	1.0	5	6
1,4-Dichlorobenzene	U	6	1.0	5	6
1,2-Dichlorobenzene	U	6	1.0	5	6
N-Butylbenzene	U	6	1.0	5	6
1,2-Dibromo-3-Chloropropane	U	6	1.0	5	6
1,2,4-Trimethylbenzene	U	6	1.0	5	6
Naphthalene	U	6	1.0	5	6
Hexachlorobutadiene	U	6	1.0	5	6
1,2,3-Trichlorobenzene	U	6	1.0	5	6
Methyl tert-butyl ether	U	12	1.0	10	12
Acetone	U	31	1.0	25	31
2-Butanone	U	31	1.0	25	31
4-methyl-2-pentanone	U	31	1.0	25	31
2-Hexanone	U	31	1.0	25	31
m+p-Xylenes	U	12	1.0	10	12
o-Xylene	U	6	1.0	5	6
Xylenes (total)	U	19	1.0	15	19
1,3,5-Trichlorobenzene	U	6	1.0	5	6
Vinyl Acetate	U	6	1.0	5	6
Carbon Disulfide	U	6	1.0	5	6
Diethyl Ether	U	6	1.0	5	6
Tetrahydrofuran	U	63	1.0	50	63
Dibromofluoromethane		71%			
1,2-Dichloroethane-D4		71%			
Toluene-D8		54%			
P-Bromofluorobenzene		* 20%			

KATAHDIN ANALYTICAL SERVICES
Report of Analytical Results

Client: Woodard & Curran
Project: Old Town
PO No:
Sample Date: 05/30/06
Received Date: 05/31/06
Extraction Date: 05/31/06
Analysis Date: 31-MAY-2006 19:28
Report Date: 06/06/2006
Matrix: SOIL
% Solids: 52.5

Lab ID: WW2637-10
Client ID: SS04
SDG: WW2637
Extracted by: KM
Extraction Method: SW846 3550
Analyst: LRS
Analysis Method: SW846 8270C
Lab Prep Batch: WG28621
Units: ug/Kg

Compound	Flags	Results	DF	PQL	Adj.PQL
Phenol	U	630	1.0	330	630
Bis(2-Chloroethyl)ether	U	630	1.0	330	630
2-Chlorophenol	U	630	1.0	330	630
1,3-Dichlorobenzene	U	630	1.0	330	630
1,4-Dichlorobenzene	U	630	1.0	330	630
1,2-Dichlorobenzene	U	630	1.0	330	630
2-Methylphenol	U	630	1.0	330	630
2,2'-Oxybis(1-Chloropropane)	U	630	1.0	330	630
N-Nitroso-di-n-propylamine	U	630	1.0	330	630
3&4-Methylphenol	U	630	1.0	330	630
Hexachloroethane	U	630	1.0	330	630
Nitrobenzene	U	630	1.0	330	630
Isophorone	U	630	1.0	330	630
2-Nitrophenol	U	630	1.0	330	630
2,4-Dimethylphenol	U	630	1.0	330	630
Bis(2-Chloroethoxy)methane	U	630	1.0	330	630
2,4-Dichlorophenol	U	630	1.0	330	630
1,2,4-Trichlorobenzene	U	630	1.0	330	630
Naphthalene	U	630	1.0	330	630
4-Chloroaniline	U	630	1.0	330	630
Hexachlorobutadiene	U	630	1.0	330	630
4-Chloro-3-Methylphenol	U	630	1.0	330	630
2-Methylnaphthalene	U	630	1.0	330	630
Hexachlorocyclopentadiene	U	630	1.0	330	630
2,4,6-Trichlorophenol	U	630	1.0	330	630
2,4,5-Trichlorophenol	U	1600	1.0	820	1600
2-Chloronaphthalene	U	630	1.0	330	630
2-Nitroaniline	U	1600	1.0	820	1600
Dimethyl Phthalate	U	630	1.0	330	630
2,6-Dinitrotoluene	U	630	1.0	330	630
Acenaphthylene	U	630	1.0	330	630
3-Nitroaniline	U	1600	1.0	820	1600
Acenaphthene	U	630	1.0	330	630
2,4-Dinitrophenol	U	1600	1.0	820	1600
Dibenzofuran	U	630	1.0	330	630
4-Nitrophenol	U	1600	1.0	820	1600
2,4-Dinitrotoluene	U	1600	1.0	820	1600
Diethylphthalate	U	630	1.0	330	630
Fluorene	U	630	1.0	330	630
4-Chlorophenyl-phenylether	U	630	1.0	330	630
4-Nitroaniline	U	1600	1.0	820	1600
4,6-Dinitro-2-Methylphenol	U	1600	1.0	820	1600
N-Nitrosodiphenylamine	U	630	1.0	330	630

KATAHDIN ANALYTICAL SERVICES
Report of Analytical Results

Client: Woodard & Curran
Project: Old Town
PO No:
Sample Date: 05/30/06
Received Date: 05/31/06
Extraction Date: 05/31/06
Analysis Date: 31-MAY-2006 19:28
Report Date: 06/06/2006
Matrix: SOIL
% Solids: 52.5

Lab ID: WW2637-10
Client ID: SS04
SDG: WW2637
Extracted by: KM
Extraction Method: SW846 3550
Analyst: LRS
Analysis Method: SW846 8270C
Lab Prep Batch: WG28621
Units: ug/Kg

Compound	Flags	Results	DF	PQL	Adj.PQL
4-Bromophenyl-phenylether	U	630	1.0	330	630
Hexachlorobenzene	U	630	1.0	330	630
Pentachlorophenol	U	1600	1.0	820	1600
Phenanthrene	U	630	1.0	330	630
Anthracene	U	630	1.0	330	630
Carbazole	U	630	1.0	330	630
Di-n-butylphthalate	U	630	1.0	330	630
Fluoranthene	U	630	1.0	330	630
Pyrene	U	630	1.0	330	630
Butylbenzylphthalate	U	630	1.0	330	630
Benzo(a)anthracene	U	630	1.0	330	630
3,3'-Dichlorobenzidine	U	630	1.0	330	630
Chrysene	U	630	1.0	330	630
bis(2-Ethylhexyl)phthalate	U	630	1.0	330	630
Di-n-octylphthalate	U	630	1.0	330	630
Benzo(b)fluoranthene	U	630	1.0	330	630
Benzo(k)fluoranthene	U	630	1.0	330	630
Benzo(a)pyrene	U	630	1.0	330	630
Indeno(1,2,3-cd)pyrene	U	630	1.0	330	630
Dibenzo(a,h)anthracene	U	630	1.0	330	630
Benzo(g,h,i)perylene	U	630	1.0	330	630
2-Fluorophenol		55%			
Phenol-D6		65%			
Nitrobenzene-D5		55%			
2-Fluorobiphenyl		63%			
2,4,6-Tribromophenol		65%			
Terphenyl-D14		87%			

Katahdin
ANALYTICAL SERVICES

REPORT OF ANALYTICAL RESULTS

Client: Erika Lloyd
Woodard & Curran
41 Hutchins Drive
Portland, ME 04102

Lab Sample ID: WW2637-010
Report Date: 6/7/2006
PO No.:
Project: Old Town

Sample Description	Matrix	Percent Solids(%)	Date Sampled	Date Received
SS04	SL	52.5	05/30/2006	05/31/2006

Parameter	Result	Units	Adjusted PQL	Dilution Factor	PQL	Analytical Method	Analysis Date	By	Prep Method	Prepped Date	By	QC	Notes
ARSENIC	U 1.	mg/Kg	1.	1	0.8	SW846 6010	6/1/06	SJC	SW846 3050	5/31/06	EAM	WE31ICS0	1
BARIUM	190.	mg/Kg	0.86	1	0.5	SW846 6010	6/1/06	SJC	SW846 3050	5/31/06	EAM	WE31ICS0	
CADMIUM	1.06	mg/Kg	0.172	5	0.02	SW846 6020	6/2/06	EAM	SW846 3050	5/31/06	EAM	WE31ICS0	
CHROMIUM	6.59	mg/Kg	0.51	5	0.06	SW846 6020	6/2/06	EAM	SW846 3050	5/31/06	EAM	WE31ICS0	
LEAD	7.70	mg/Kg	0.17	5	0.02	SW846 6020	6/2/06	EAM	SW846 3050	5/31/06	EAM	WE31ICS0	
MERCURY	U 0.067	ug/g	0.067	1	0.04	SW846 7471	6/2/06	MW	SW846 7471	6/1/06	MW	WF01HGS0	1
SELENIUM	U 1.7	mg/Kg	1.7	1	1	SW846 6010	6/1/06	SJC	SW846 3050	5/31/06	EAM	WE31ICS0	1
SILVER	U 2.6	mg/Kg	2.6	1	1.5	SW846 6010	6/1/06	SJC	SW846 3050	5/31/06	EAM	WE31ICS0	1

1 The laboratory's Practical Quantitation Level could not be achieved for this parameter due to sample composition, matrix effects,sample volume, or quantity used for analysis.

Report of Analytical Results

Client: Erika Lloyd
Woodard & Curran
41 Hutchins Drive
Portland,ME 04102

Lab Sample ID: WW2637-10
Report Date: 07-JUN-06
Client PO:
Project: Old Town
SDG: WW2637

Sample Description	Matrix	Date Sampled	Date Received
SS04	SL	30-MAY-06	31-MAY-06

Parameter	Result	Adj PQL	Anal. Method	QC.Batch	Anal. Date	Prep. Method	Prep. Date	Analyst	Footnotes
Total Solids	52. %	1	CLP SOW 788	WG28725	02-JUN-06 08:46:00	CLP SOW 788	01-JUN-06	JF	

KATAHDIN ANALYTICAL SERVICES
Report of Analytical Results

Client: Woodard & Curran
Project: Old Town
PO No:
Sample Date: 05/30/06
Received Date: 05/31/06
Extraction Date:
Analysis Date: 31-MAY-2006 13:47
Report Date: 06/02/2006
Matrix: SOIL
% Solids: 60.0

Lab ID: WW2637-11
Client ID: SS05
SDG: WW2637
Extracted by:
Extraction Method: SW846 5035
Analyst: ALH
Analysis Method: SW846 8260B
Lab Prep Batch: WG28636
Units: ug/Kg

Compound	Flags	Results	DF	PQL	Adj.PQL
Dichlorodifluoromethane	U	10	1.0	10	10
Chloromethane	U	10	1.0	10	10
Vinyl chloride	U	10	1.0	10	10
Bromomethane	U	10	1.0	10	10
Chloroethane	U	10	1.0	10	10
Trichlorofluoromethane	U	10	1.0	10	10
1,1-Dichloroethene	U	5	1.0	5	5
Methylene Chloride	U	25	1.0	25	25
trans-1,2-Dichloroethene	U	5	1.0	5	5
1,1-Dichloroethane	U	5	1.0	5	5
cis-1,2-Dichloroethene	U	5	1.0	5	5
1,2-Dichloroethylene (total)	U	10	1.0	10	10
2,2-Dichloropropane	U	5	1.0	5	5
Chloroform	U	5	1.0	5	5
Bromochloromethane	U	5	1.0	5	5
1,1,1-Trichloroethane	U	5	1.0	5	5
1,2-Dichloroethane	U	5	1.0	5	5
1,1-Dichloropropene	U	5	1.0	5	5
Carbon Tetrachloride	U	5	1.0	5	5
Benzene	U	5	1.0	5	5
1,2-Dichloropropane	U	5	1.0	5	5
Trichloroethene	U	5	1.0	5	5
Dibromomethane	U	5	1.0	5	5
Bromodichloromethane	U	5	1.0	5	5
cis-1,3-dichloropropene	U	5	1.0	5	5
Toluene	U	5	1.0	5	5
trans-1,3-Dichloropropene	U	5	1.0	5	5
1,1,2-Trichloroethane	U	5	1.0	5	5
1,3-Dichloropropane	U	5	1.0	5	5
Dibromochloromethane	U	5	1.0	5	5
Tetrachloroethene	U	5	1.0	5	5
1,2-Dibromoethane	U	5	1.0	5	5
Chlorobenzene	U	5	1.0	5	5
1,1,1,2-Tetrachloroethane	U	5	1.0	5	5
Ethylbenzene	U	5	1.0	5	5
Bromoform	U	5	1.0	5	5
Styrene	U	5	1.0	5	5
1,1,2,2-Tetrachloroethane	U	5	1.0	5	5
1,2,3-Trichloropropane	U	5	1.0	5	5
Isopropylbenzene	U	5	1.0	5	5
Bromobenzene	U	5	1.0	5	5
2-Chlorotoluene	U	5	1.0	5	5
N-Propylbenzene	U	5	1.0	5	5

KATAHDIN ANALYTICAL SERVICES
Report of Analytical Results

Client: Woodard & Curran
Project: Old Town
PO No:
Sample Date: 05/30/06
Received Date: 05/31/06
Extraction Date:
Analysis Date: 31-MAY-2006 13:47
Report Date: 06/02/2006
Matrix: SOIL
% Solids: 60.0

Lab ID: WW2637-11
Client ID: SS05
SDG: WW2637
Extracted by:
Extraction Method: SW846 5035
Analyst: ALH
Analysis Method: SW846 8260B
Lab Prep Batch: WG28636
Units: ug/Kg

Compound	Flags	Results	DF	PQL	Adj.PQL
4-Chlorotoluene	U	5	1.0	5	5
1,3,5-Trimethylbenzene	U	5	1.0	5	5
tert-Butylbenzene	U	5	1.0	5	5
1,2,4-Trichlorobenzene	U	5	1.0	5	5
sec-Butylbenzene	U	5	1.0	5	5
1,3-Dichlorobenzene	U	5	1.0	5	5
P-Isopropyltoluene	U	5	1.0	5	5
1,4-Dichlorobenzene	U	5	1.0	5	5
1,2-Dichlorobenzene	U	5	1.0	5	5
N-Butylbenzene	U	5	1.0	5	5
1,2-Dibromo-3-Chloropropane	U	5	1.0	5	5
1,2,4-Trimethylbenzene	U	5	1.0	5	5
Naphthalene	U	5	1.0	5	5
Hexachlorobutadiene	U	5	1.0	5	5
1,2,3-Trichlorobenzene	U	5	1.0	5	5
Methyl tert-butyl ether	U	10	1.0	10	10
Acetone		71	1.0	25	25
2-Butanone	U	25	1.0	25	25
4-methyl-2-pentanone	U	25	1.0	25	25
2-Hexanone	U	25	1.0	25	25
m+p-Xylenes	U	10	1.0	10	10
o-Xylene	U	5	1.0	5	5
Xylenes (total)	U	15	1.0	15	15
1,3,5-Trichlorobenzene	U	5	1.0	5	5
Vinyl Acetate	U	5	1.0	5	5
Carbon Disulfide	U	5	1.0	5	5
Diethyl Ether	U	5	1.0	5	5
Tetrahydrofuran	U	50	1.0	50	50
Dibromofluoromethane		60%			
1,2-Dichloroethane-D4		63%			
Toluene-D8		50%			
P-Bromofluorobenzene		* 22%			

KATAHDIN ANALYTICAL SERVICES
Report of Analytical Results

Client: Woodard & Curran
Project: Old Town
PO No:
Sample Date: 05/30/06
Received Date: 05/31/06
Extraction Date: 05/31/06
Analysis Date: 02-JUN-2006 09:44
Report Date: 06/06/2006
Matrix: SOIL
% Solids: 60.0

Lab ID: WW2637-11RA
Client ID: SS05
SDG: WW2637
Extracted by: KM
Extraction Method: SW846 3550
Analyst: LRS
Analysis Method: SW846 8270C
Lab Prep Batch: WG28621
Units: ug/Kg

Compound	Flags	Results	DF	PQL	Adj.PQL
Phenol	U	550	1.0	330	550
Bis(2-Chloroethyl)ether	U	550	1.0	330	550
2-Chlorophenol	U	550	1.0	330	550
1,3-Dichlorobenzene	U	550	1.0	330	550
1,4-Dichlorobenzene	U	550	1.0	330	550
1,2-Dichlorobenzene	U	550	1.0	330	550
2-Methylphenol	U	550	1.0	330	550
2,2'-Oxybis(1-Chloropropane)	U	550	1.0	330	550
N-Nitroso-di-n-propylamine	U	550	1.0	330	550
3&4-Methylphenol	U	550	1.0	330	550
Hexachloroethane	U	550	1.0	330	550
Nitrobenzene	U	550	1.0	330	550
Isophorone	U	550	1.0	330	550
2-Nitrophenol	U	550	1.0	330	550
2,4-Dimethylphenol	U	550	1.0	330	550
Bis(2-Chloroethoxy)methane	U	550	1.0	330	550
2,4-Dichlorophenol	U	550	1.0	330	550
1,2,4-Trichlorobenzene	U	550	1.0	330	550
Naphthalene	U	550	1.0	330	550
4-Chloroaniline	U	550	1.0	330	550
Hexachlorobutadiene	U	550	1.0	330	550
4-Chloro-3-Methylphenol	U	550	1.0	330	550
2-Methylnaphthalene	U	550	1.0	330	550
Hexachlorocyclopentadiene	U	550	1.0	330	550
2,4,6-Trichlorophenol	U	550	1.0	330	550
2,4,5-Trichlorophenol	U	1400	1.0	820	1400
2-Chloronaphthalene	U	550	1.0	330	550
2-Nitroaniline	U	1400	1.0	820	1400
Dimethyl Phthalate	U	550	1.0	330	550
2,6-Dinitrotoluene	U	550	1.0	330	550
Acenaphthylene	U	550	1.0	330	550
3-Nitroaniline	U	1400	1.0	820	1400
Acenaphthene	U	550	1.0	330	550
2,4-Dinitrophenol	U	1400	1.0	820	1400
Dibenzofuran	U	550	1.0	330	550
4-Nitrophenol	U	1400	1.0	820	1400
2,4-Dinitrotoluene	U	1400	1.0	820	1400
Diethylphthalate	U	550	1.0	330	550
Fluorene	U	550	1.0	330	550
4-Chlorophenyl-phenylether	U	550	1.0	330	550
4-Nitroaniline	U	1400	1.0	820	1400
4,6-Dinitro-2-Methylphenol	U	1400	1.0	820	1400
N-Nitrosodiphenylamine	U	550	1.0	330	550

KATAHDIN ANALYTICAL SERVICES
Report of Analytical Results

Client: Woodard & Curran
Project: Old Town
PO No:
Sample Date: 05/30/06
Received Date: 05/31/06
Extraction Date: 05/31/06
Analysis Date: 02-JUN-2006 09:44
Report Date: 06/06/2006
Matrix: SOIL
% Solids: 60.0

Lab ID: WW2637-11RA
Client ID: SS05
SDG: WW2637
Extracted by: KM
Extraction Method: SW846 3550
Analyst: LRS
Analysis Method: SW846 8270C
Lab Prep Batch: WG28621
Units: ug/Kg

Compound	Flags	Results	DF	PQL	Adj.PQL
4-Bromophenyl-phenylether	U	550	1.0	330	550
Hexachlorobenzene	U	550	1.0	330	550
Pentachlorophenol	U	1400	1.0	820	1400
Phenanthrene	U	550	1.0	330	550
Anthracene	U	550	1.0	330	550
Carbazole	U	550	1.0	330	550
Di-n-butylphthalate	U	550	1.0	330	550
Fluoranthene	U	550	1.0	330	550
Pyrene	U	550	1.0	330	550
Butylbenzylphthalate	U	550	1.0	330	550
Benzo(a)anthracene	U	550	1.0	330	550
3,3'-Dichlorobenzidine	U	550	1.0	330	550
Chrysene	U	550	1.0	330	550
bis(2-Ethylhexyl)phthalate	U	550	1.0	330	550
Di-n-octylphthalate	U	550	1.0	330	550
Benzo(b)fluoranthene	U	550	1.0	330	550
Benzo(k)fluoranthene	U	550	1.0	330	550
Benzo(a)pyrene	U	550	1.0	330	550
Indeno(1,2,3-cd)pyrene	U	550	1.0	330	550
Dibenzo(a,h)anthracene	U	550	1.0	330	550
Benzo(g,h,i)perylene	U	550	1.0	330	550
2-Fluorophenol		52%			
Phenol-D6		61%			
Nitrobenzene-D5		54%			
2-Fluorobiphenyl		61%			
2,4,6-Tribromophenol		63%			
Terphenyl-D14		75%			

Katahdin ANALYTICAL SERVICES

REPORT OF ANALYTICAL RESULTS

Client: Erika Lloyd
Woodard & Curran
41 Hutchins Drive
Portland, ME 04102

Lab Sample ID: WW2637-011
Report Date: 6/7/2006
PO No.:
Project: Old Town

Sample Description	Matrix	Percent Solids(%)	Date Sampled	Date Received
SS05	SL	60.0	05/30/2006	05/31/2006

Parameter	Result	Units	Adjusted PQL	Dilution Factor	PQL	Analytical Method	Analysis Date	By	Prep Method	Prepped Date	By	QC	Notes
ARSENIC	1.	mg/Kg	1.	1	0.8	SW846 6010	6/1/06	SJC	SW846 3050	5/31/06	EAM	WE31ICS0	
BARIUM	85.6	mg/Kg	0.74	1	0.5	SW846 6010	6/1/06	SJC	SW846 3050	5/31/06	EAM	WE31ICS0	
CADMIUM	0.393	mg/Kg	0.149	5	0.02	SW846 6020	6/2/06	EAM	SW846 3050	5/31/06	EAM	WE31ICS0	
CHROMIUM	29.8	mg/Kg	0.45	5	0.06	SW846 6020	6/2/06	EAM	SW846 3050	5/31/06	EAM	WE31ICS0	
LEAD	4.43	mg/Kg	0.15	5	0.02	SW846 6020	6/2/06	EAM	SW846 3050	5/31/06	EAM	WE31ICS0	
MERCURY	U 0.058	ug/g	0.058	1	0.04	SW846 7471	6/2/06	MW	SW846 7471	6/1/06	MW	WF01HGS0	1
SELENIUM	U 1.5	mg/Kg	1.5	1	1	SW846 6010	6/1/06	SJC	SW846 3050	5/31/06	EAM	WE31ICS0	1
SILVER	U 2.2	mg/Kg	2.2	1	1.5	SW846 6010	6/1/06	SJC	SW846 3050	5/31/06	EAM	WE31ICS0	1

1 The laboratory's Practical Quantitation Level could not be achieved for this parameter due to sample composition, matrix effects,sample volume, or quantity used for analysis.

Report of Analytical Results

Client: Erika Lloyd
Woodard & Curran
41 Hutchins Drive
Portland,ME 04102

Lab Sample ID: WW2637-11
Report Date: 07-JUN-06
Client PO:
Project: Old Town
SDG: WW2637

Sample Description	Matrix	Date Sampled	Date Received
SS05	SL	30-MAY-06	31-MAY-06

Parameter	Result	Adj PQL	Anal. Method	QC.Batch	Anal. Date	Prep. Method	Prep. Date	Analyst	Footnotes
Total Solids	60. %	1	CLP SOW 788	WG28725	02-JUN-06 08:47:00	CLP SOW 788	01-JUN-06	JF	

KATAHDIN ANALYTICAL SERVICES
Report of Analytical Results

Client: Woodard & Curran
Project: Old Town
PO No:
Sample Date: 05/30/06
Received Date: 05/31/06
Extraction Date: 05/31/06
Analysis Date: 01-JUN-2006 14:54
Report Date: 06/02/2006
Matrix: SOIL
% Solids: 36.0

Lab ID: WW2637-12
Client ID: SS07
SDG: WW2637
Extracted by: KM
Extraction Method: SW846 3550
Analyst: JLP
Analysis Method: SW846 8082
Lab Prep Batch: WG28624
Units: ug/Kg

Compound	Flags	Results	DF	PQL	Adj.PQL
Aroclor-1016	U	47	1.0	17	47
Aroclor-1221	U	47	1.0	17	47
Aroclor-1232	U	47	1.0	17	47
Aroclor-1242	U	47	1.0	17	47
Aroclor-1248	U	47	1.0	17	47
Aroclor-1254	U	47	1.0	17	47
Aroclor-1260	U	47	1.0	17	47
Tetrachloro-m-xylene		79%			
Decachlorobiphenyl		79%			

Report of Analytical Results

Client: Erika Lloyd
Woodard & Curran
41 Hutchins Drive
Portland,ME 04102

Lab Sample ID: WW2637-12
Report Date: 07-JUN-06
Client PO:
Project: Old Town
SDG: WW2637

Sample Description	Matrix	Date Sampled	Date Received
SS07	SL	30-MAY-06	31-MAY-06

Parameter	Result	Adj PQL	Anal. Method	QC.Batch	Anal. Date	Prep. Method	Prep. Date	Analyst	Footnotes
Total Solids	36. %	1	CLP SOW 788	WG28725	02-JUN-06 08:39:00	CLP SOW 788	01-JUN-06	JF	

KATAHDIN ANALYTICAL SERVICES
Report of Analytical Results

Client: Woodard & Curran
Project: Old Town
PO No:
Sample Date: 05/30/06
Received Date: 05/31/06
Extraction Date: 05/31/06
Analysis Date: 01-JUN-2006 15:22
Report Date: 06/02/2006
Matrix: SOIL
% Solids: 35.8

Lab ID: WW2637-13
Client ID: SS08
SDG: WW2637
Extracted by: KM
Extraction Method: SW846 3550
Analyst: JLP
Analysis Method: SW846 8082
Lab Prep Batch: WG28624
Units: ug/Kg

Compound	Flags	Results	DF	PQL	Adj.PQL
Aroclor-1016	U	47	1.0	17	47
Aroclor-1221	U	47	1.0	17	47
Aroclor-1232	U	47	1.0	17	47
Aroclor-1242	U	47	1.0	17	47
Aroclor-1248	U	47	1.0	17	47
Aroclor-1254	U	47	1.0	17	47
Aroclor-1260	U	47	1.0	17	47
Tetrachloro-m-xylene		67%			
Decachlorobiphenyl		64%			

Report of Analytical Results

Client: Erika Lloyd
Woodard & Curran
41 Hutchins Drive
Portland,ME 04102

Lab Sample ID: WW2637-13
Report Date: 07-JUN-06
Client PO:
Project: Old Town
SDG: WW2637

Sample Description	Matrix	Date Sampled	Date Received
SS08	SL	30-MAY-06	31-MAY-06

Parameter	Result	Adj PQL	Anal. Method	QC.Batch	Anal. Date	Prep. Method	Prep. Date	Analyst	Footnotes
Total Solids	36. %	1	CLP SOW 788	WG28725	02-JUN-06 08:39:00	CLP SOW 788	01-JUN-06	JF	

KATAHDIN ANALYTICAL SERVICES
Report of Analytical Results

Client: Woodard & Curran
Project: Old Town
PO No:
Sample Date: 05/30/06
Received Date: 05/31/06
Extraction Date:
Analysis Date: 31-MAY-2006 14:23
Report Date: 06/07/2006
Matrix: SOIL
% Solids: 100

Lab ID: WW2637-14
Client ID: SS81
SDG: WW2637
Extracted by:
Extraction Method: SW846 5035
Analyst: ALH
Analysis Method: SW846 8260B
Lab Prep Batch: WG28636
Units: ug/Kg

Compound	Flags	Results	DF	PQL	Adj.PQL
Dichlorodifluoromethane	U	10	1.0	10	10
Chloromethane	U	10	1.0	10	10
Vinyl chloride	U	10	1.0	10	10
Bromomethane	U	10	1.0	10	10
Chloroethane	U	10	1.0	10	10
Trichlorofluoromethane	U	10	1.0	10	10
1,1-Dichloroethene	U	5	1.0	5	5
Methylene Chloride	U	25	1.0	25	25
trans-1,2-Dichloroethene	U	5	1.0	5	5
1,1-Dichloroethane	U	5	1.0	5	5
cis-1,2-Dichloroethene	U	5	1.0	5	5
1,2-Dichloroethylene (total)	U	10	1.0	10	10
2,2-Dichloropropane	U	5	1.0	5	5
Chloroform	U	5	1.0	5	5
Bromochloromethane	U	5	1.0	5	5
1,1,1-Trichloroethane	U	5	1.0	5	5
1,2-Dichloroethane	U	5	1.0	5	5
1,1-Dichloropropene	U	5	1.0	5	5
Carbon Tetrachloride	U	5	1.0	5	5
Benzene	U	5	1.0	5	5
1,2-Dichloropropane	U	5	1.0	5	5
Trichloroethene	U	5	1.0	5	5
Dibromomethane	U	5	1.0	5	5
Bromodichloromethane	U	5	1.0	5	5
cis-1,3-dichloropropene	U	5	1.0	5	5
Toluene	U	5	1.0	5	5
trans-1,3-Dichloropropene	U	5	1.0	5	5
1,1,2-Trichloroethane	U	5	1.0	5	5
1,3-Dichloropropane	U	5	1.0	5	5
Dibromochloromethane	U	5	1.0	5	5
Tetrachloroethene	U	5	1.0	5	5
1,2-Dibromoethane	U	5	1.0	5	5
Chlorobenzene	U	5	1.0	5	5
1,1,1,2-Tetrachloroethane	U	5	1.0	5	5
Ethylbenzene	U	5	1.0	5	5
Bromoform	U	5	1.0	5	5
Styrene	U	5	1.0	5	5
1,1,2,2-Tetrachloroethane	U	5	1.0	5	5
1,2,3-Trichloropropane	U	5	1.0	5	5
Isopropylbenzene	U	5	1.0	5	5
Bromobenzene	U	5	1.0	5	5
2-Chlorotoluene	U	5	1.0	5	5
N-Propylbenzene	U	5	1.0	5	5

KATAHDIN ANALYTICAL SERVICES
Report of Analytical Results

Client: Woodard & Curran
Project: Old Town
PO No:
Sample Date: 05/30/06
Received Date: 05/31/06
Extraction Date:
Analysis Date: 31-MAY-2006 14:23
Report Date: 06/07/2006
Matrix: SOIL
% Solids: 100

Lab ID: WW2637-14
Client ID: SS81
SDG: WW2637
Extracted by:
Extraction Method: SW846 5035
Analyst: ALH
Analysis Method: SW846 8260B
Lab Prep Batch: WG28636
Units: ug/Kg

Compound	Flags	Results	DF	PQL	Adj.PQL
4-Chlorotoluene	U	5	1.0	5	5
1,3,5-Trimethylbenzene	U	5	1.0	5	5
tert-Butylbenzene	U	5	1.0	5	5
1,2,4-Trichlorobenzene	U	5	1.0	5	5
sec-Butylbenzene	U	5	1.0	5	5
1,3-Dichlorobenzene	U	5	1.0	5	5
P-Isopropyltoluene	U	5	1.0	5	5
1,4-Dichlorobenzene	U	5	1.0	5	5
1,2-Dichlorobenzene	U	5	1.0	5	5
N-Butylbenzene	U	5	1.0	5	5
1,2-Dibromo-3-Chloropropane	U	5	1.0	5	5
1,2,4-Trimethylbenzene	U	5	1.0	5	5
Naphthalene	U	5	1.0	5	5
Hexachlorobutadiene	U	5	1.0	5	5
1,2,3-Trichlorobenzene	U	5	1.0	5	5
Methyl tert-butyl ether	U	10	1.0	10	10
Acetone		70	1.0	25	25
2-Butanone	U	25	1.0	25	25
4-methyl-2-pentanone	U	25	1.0	25	25
2-Hexanone	U	25	1.0	25	25
m+p-Xylenes	U	10	1.0	10	10
o-Xylene	U	5	1.0	5	5
Xylenes (total)	U	15	1.0	15	15
1,3,5-Trichlorobenzene	U	5	1.0	5	5
Vinyl Acetate	U	5	1.0	5	5
Carbon Disulfide	U	5	1.0	5	5
Diethyl Ether	U	5	1.0	5	5
Tetrahydrofuran	U	50	1.0	50	50
Dibromofluoromethane		70%			
1,2-Dichloroethane-D4		69%			
Toluene-D8		60%			
P-Bromofluorobenzene		* 32%			

KATAHDIN ANALYTICAL SERVICES
Report of Analytical Results

Client: Woodard & Curran
Project: Old Town
PO No:
Sample Date: 05/09/06
Received Date: 05/31/06
Extraction Date:
Analysis Date: 31-MAY-2006 12:33
Report Date: 06/02/2006
Matrix: WATER
% Solids: NA

Lab ID: WW2637-15
Client ID: TRIP BLANK
SDG: WW2637
Extracted by:
Extraction Method: SW846 5030
Analyst: SKT
Analysis Method: SW846 8260B
Lab Prep Batch: WG28646
Units: ug/l

Compound	Flags	Results	DF	PQL	Adj.PQL
Dichlorodifluoromethane	U	10	1.0	10	10
Chloromethane	U	10	1.0	10	10
Vinyl chloride	U	10	1.0	10	10
Bromomethane	U	10	1.0	10	10
Chloroethane	U	10	1.0	10	10
Trichlorofluoromethane	U	10	1.0	10	10
1,1-Dichloroethene	U	5	1.0	5	5
Methylene Chloride	U	5	1.0	5	5
trans-1,2-Dichloroethene	U	5	1.0	5	5
1,1-Dichloroethane	U	5	1.0	5	5
cis-1,2-Dichloroethene	U	5	1.0	5	5
1,2-Dichloroethylene (total)	U	10	1.0	10	10
2,2-Dichloropropane	U	5	1.0	5	5
Chloroform	U	5	1.0	5	5
Bromochloromethane	U	5	1.0	5	5
1,1,1-Trichloroethane	U	5	1.0	5	5
1,2-Dichloroethane	U	5	1.0	5	5
1,1-Dichloropropene	U	5	1.0	5	5
Carbon Tetrachloride	U	5	1.0	5	5
Benzene	U	5	1.0	5	5
1,2-Dichloropropane	U	5	1.0	5	5
Trichloroethene	U	5	1.0	5	5
Dibromomethane	U	5	1.0	5	5
Bromodichloromethane	U	5	1.0	5	5
cis-1,3-dichloropropene	U	5	1.0	5	5
Toluene	U	5	1.0	5	5
trans-1,3-Dichloropropene	U	5	1.0	5	5
1,1,2-Trichloroethane	U	5	1.0	5	5
1,3-Dichloropropane	U	5	1.0	5	5
Dibromochloromethane	U	5	1.0	5	5
Tetrachloroethene	U	5	1.0	5	5
1,2-Dibromoethane	U	5	1.0	5	5
Chlorobenzene	U	5	1.0	5	5
1,1,1,2-Tetrachloroethane	U	5	1.0	5	5
Ethylbenzene	U	5	1.0	5	5
Bromoform	U	5	1.0	5	5
Styrene	U	5	1.0	5	5
1,1,2,2-Tetrachloroethane	U	5	1.0	5	5
1,2,3-Trichloropropane	U	5	1.0	5	5
Isopropylbenzene	U	5	1.0	5	5
Bromobenzene	U	5	1.0	5	5
2-Chlorotoluene	U	5	1.0	5	5
N-Propylbenzene	U	5	1.0	5	5

KATAHDIN ANALYTICAL SERVICES
Report of Analytical Results

Client: Woodard & Curran
Project: Old Town
PO No:
Sample Date: 05/09/06
Received Date: 05/31/06
Extraction Date:
Analysis Date: 31-MAY-2006 12:33
Report Date: 06/02/2006
Matrix: WATER
% Solids: NA

Lab ID: WW2637-15
Client ID: TRIP BLANK
SDG: WW2637
Extracted by:
Extraction Method: SW846 5030
Analyst: SKT
Analysis Method: SW846 8260B
Lab Prep Batch: WG28646
Units: ug/l

Compound	Flags	Results	DF	PQL	Adj.PQL
4-Chlorotoluene	U	5	1.0	5	5
1,3,5-Trimethylbenzene	U	5	1.0	5	5
tert-Butylbenzene	U	5	1.0	5	5
1,2,4-Trichlorobenzene	U	5	1.0	5	5
sec-Butylbenzene	U	5	1.0	5	5
1,3-Dichlorobenzene	U	5	1.0	5	5
P-Isopropyltoluene	U	5	1.0	5	5
1,4-Dichlorobenzene	U	5	1.0	5	5
1,2-Dichlorobenzene	U	5	1.0	5	5
N-Butylbenzene	U	5	1.0	5	5
1,2-Dibromo-3-Chloropropane	U	5	1.0	5	5
1,2,4-Trimethylbenzene	U	5	1.0	5	5
Naphthalene	U	5	1.0	5	5
Hexachlorobutadiene	U	5	1.0	5	5
1,2,3-Trichlorobenzene	U	5	1.0	5	5
Methyl tert-butyl ether	U	10	1.0	10	10
Acetone	U	25	1.0	25	25
2-Butanone	U	25	1.0	25	25
4-methyl-2-pentanone	U	25	1.0	25	25
2-Hexanone	U	25	1.0	25	25
m+p-Xylenes	U	10	1.0	10	10
o-Xylene	U	5	1.0	5	5
Xylenes (total)	U	15	1.0	15	15
1,3,5-Trichlorobenzene	U	5	1.0	5	5
Vinyl Acetate	U	5	1.0	5	5
Carbon Disulfide	U	5	1.0	5	5
Diethyl Ether	U	5	1.0	5	5
Tetrahydrofuran	U	25	1.0	25	25
Dibromofluoromethane		87%			
1,2-Dichloroethane-D4		90%			
Toluene-D8		85%			
P-Bromofluorobenzene		84%			

FORM 4
VOLATILE METHOD BLANK SUMMARY

CLIENT SAMPLE ID

| WG28646-BLANK |

Lab Name: KATAHDIN ANALYTICAL SERVICES Lab Code: KAS

Project: OLD TOWN SDG No.: WW2637

Lab File ID: T1134 Lab Sample ID: WG28646-2

Date Analyzed: 05/31/06 Time Analyzed: 1048

GC Column: RTX-VMS ID: 0.18 (mm) Heated Purge: (Y/N) N

Instrument ID: GCMS-T

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES, MS and MSD:

	CLIENT SAMPLE ID	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED	TIME ANALYZED
01	WG28646-LCS	WG28646-1	T1132	05/31/06	0935
02	TRIP BLANK	WW2637-15	T1137	05/31/06	1233
03	SB-25	WW2637-1	T1143	05/31/06	1537
04	MW-1	WW2637-2	T1144	05/31/06	1608
05	MW-2	WW2637-3	T1145	05/31/06	1638
06	SB-26	WW2637-4	T1146	05/31/06	1709
07	SB-100	WW2637-5	T1147	05/31/06	1740
08	SB-25MS	WG28646-3	T1148	05/31/06	1811
09	SB-25MSD	WG28646-4	T1149	05/31/06	1842
10					
11					
12					
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28					
29					
30					

COMMENTS:

FORM IV VOA

KATAHDIN ANALYTICAL SERVICES
Report of Analytical Results

Client:
Project: Old Town
PO No:
Sample Date:
Received Date:
Extraction Date:
Analysis Date: 31-MAY-2006 10:48
Report Date: 06/07/2006
Matrix: WATER
% Solids: NA

Lab ID: WG28646-2
Client ID: WG28646-Blank
SDG: WW2637
Extracted by:
Extraction Method: SW846 5030
Analyst: SKT
Analysis Method: SW846 8260B
Lab Prep Batch: WG28646
Units: ug/l

Compound	Flags	Results	DF	PQL	Adj.PQL
Dichlorodifluoromethane	U	1	1.0	1	1
Chloromethane	U	1	1.0	1	1
Vinyl chloride	U	1	1.0	1	1
Bromomethane	U	1	1.0	1	1
Chloroethane	U	1	1.0	1	1
Trichlorofluoromethane	U	1	1.0	1	1
1,1-Dichloroethene	U	1	1.0	1	1
Methylene Chloride	U	5	1.0	5	5
trans-1,2-Dichloroethene	U	1	1.0	1	1
1,1-Dichloroethane	U	1	1.0	1	1
cis-1,2-Dichloroethene	U	1	1.0	1	1
1,2-Dichloroethylene (total)	U	2	1.0	2	2
2,2-Dichloropropane	U	1	1.0	1	1
Chloroform	U	1	1.0	1	1
Bromochloromethane	U	1	1.0	1	1
1,1,1-Trichloroethane	U	1	1.0	1	1
1,2-Dichloroethane	U	1	1.0	1	1
1,1-Dichloropropene	U	1	1.0	1	1
Carbon Tetrachloride	U	1	1.0	1	1
Benzene	U	1	1.0	1	1
1,2-Dichloropropane	U	1	1.0	1	1
Trichloroethene	U	1	1.0	1	1
Dibromomethane	U	1	1.0	1	1
Bromodichloromethane	U	1	1.0	1	1
cis-1,3-dichloropropene	U	1	1.0	1	1
Toluene	U	1	1.0	1	1
trans-1,3-Dichloropropene	U	1	1.0	1	1
1,1,2-Trichloroethane	U	1	1.0	1	1
1,3-Dichloropropane	U	1	1.0	1	1
Dibromochloromethane	U	1	1.0	1	1
Tetrachloroethene	U	1	1.0	1	1
1,2-Dibromoethane	U	1	1.0	1	1
Chlorobenzene	U	1	1.0	1	1
1,1,1,2-Tetrachloroethane	U	1	1.0	1	1
Ethylbenzene	U	1	1.0	1	1
Bromoform	U	1	1.0	1	1
Styrene	U	1	1.0	1	1
1,1,2,2-Tetrachloroethane	U	1	1.0	1	1
1,2,3-Trichloropropane	U	1	1.0	1	1
Isopropylbenzene	U	1	1.0	1	1
Bromobenzene	U	1	1.0	1	1
2-Chlorotoluene	U	1	1.0	1	1
N-Propylbenzene	U	1	1.0	1	1

KATAHDIN ANALYTICAL SERVICES
Report of Analytical Results

Client:
Project: Old Town
PO No:
Sample Date:
Received Date:
Extraction Date:
Analysis Date: 31-MAY-2006 10:48
Report Date: 06/07/2006
Matrix: WATER
% Solids: NA

Lab ID: WG28646-2
Client ID: WG28646-Blank
SDG: WW2637
Extracted by:
Extraction Method: SW846 5030
Analyst: SKT
Analysis Method: SW846 8260B
Lab Prep Batch: WG28646
Units: ug/l

Compound	Flags	Results	DF	PQL	Adj.PQL
4-Chlorotoluene	U	1	1.0	1	1
1,3,5-Trimethylbenzene	U	1	1.0	1	1
tert-Butylbenzene	U	1	1.0	1	1
1,2,4-Trichlorobenzene	U	1	1.0	1	1
sec-Butylbenzene	U	1	1.0	1	1
1,3-Dichlorobenzene	U	1	1.0	1	1
P-Isopropyltoluene	U	1	1.0	1	1
1,4-Dichlorobenzene	U	1	1.0	1	1
1,2-Dichlorobenzene	U	1	1.0	1	1
N-Butylbenzene	U	1	1.0	1	1
1,2-Dibromo-3-Chloropropane	U	1	1.0	1	1
1,2,4-Trimethylbenzene	U	1	1.0	1	1
Naphthalene		5	1.0	1	1
Hexachlorobutadiene	U	1	1.0	1	1
1,2,3-Trichlorobenzene	U	1	1.0	1	1
Methyl tert-butyl ether	U	1	1.0	1	1
Acetone	U	5	1.0	5	5
2-Butanone	U	5	1.0	5	5
4-methyl-2-pentanone	U	5	1.0	5	5
2-Hexanone	U	5	1.0	5	5
m+p-Xylenes	U	2	1.0	2	2
o-Xylene	U	1	1.0	1	1
Xylenes (total)	U	3	1.0	3	3
1,3,5-Trichlorobenzene	U	1	1.0	1	1
Vinyl Acetate	U	1	1.0	1	1
Carbon Disulfide	U	1	1.0	1	1
Diethyl Ether	U	1	1.0	1	1
Tetrahydrofuran	U	5	1.0	5	5
Dibromofluoromethane		83%			
1,2-Dichloroethane-D4		84%			
Toluene-D8		84%			
P-Bromofluorobenzene		84%			

KATAHDIN ANALYTICAL SERVICES
LAB CONTROL SAMPLE

Client:
Project: Old Town
PO No:
Sample Date:
Received Date:
Extraction Date:
Analysis Date: 05/31/06
Report Date: 06/07/2006
Matrix: WATER

Lab ID: WG28646-1
Client ID: WG28646-LCS
SDG: WW2637
Extracted by:
Extraction Method: SW846 5030
Analyst: SKT
Analysis Method: SW846 8260B
Lab Prep Batch: WG28646
Units: ug/l

COMPOUND	LCS SPIKE	SAMPLE CONC.	LCS CONC.	%REC.	QC. LIMITS
Dichlorodifluoromethane	50	NA	30	61	0-200
Chloromethane	50	NA	35	71	40-125
Vinyl chloride	50	NA	40	80	27-143
Bromomethane	50	NA	46	93	52-141
Chloroethane	50	NA	49	99	64-124
Trichlorofluoromethane	50	NA	63	126	63-153
Diethyl Ether	50	NA	50	101	63-146
Tertiary-butyl alcohol	250	NA	290	116	50-150
1,1-Dichloroethene	50	NA	47	94	85-129
Carbon Disulfide	50	NA	51	102	80-134
Freon-113	50	NA	53	106	79-129
Iodomethane	50	NA	48	96	67-144
Acrolein	250	NA	234	94	0-200
Methylene Chloride	50	NA	48	96	86-124
Acetone	50	NA	64	129	41-180
Isobutyl Alcohol	1000	NA	1120	112	20-176
trans-1,2-Dichloroethene	50	NA	46	91	80-128
Allyl Chloride	50	NA	51	102	68-133
Methyl tert-butyl ether	100	NA	107	107	79-132
Acetonitrile	500	NA	458	92	35-171
Di-isopropyl ether	50	NA	49	98	82-122
Chloroprene	50	NA	51	102	77-133
Methacrylonitrile	500	NA	484	97	73-137
Propionitrile	500	NA	485	97	54-148
1,1-Dichloroethane	50	NA	48	97	90-127
Acrylonitrile	250	NA	238	95	70-135
Ethyl tertiary-butyl ether	50	NA	53	105	84-122
Vinyl Acetate	50	NA	45	90	47-122
cis-1,2-Dichloroethene	50	NA	44	89	84-121
1,2-Dichloroethylene (total)	100	NA	90	90	83-123
Methyl Methacrylate	50	NA	52	104	73-133
2,2-Dichloropropane	50	NA	44	89	68-149
Bromochloromethane	50	NA	48	97	87-120
Chloroform	50	NA	49	99	87-127
Carbon Tetrachloride	50	NA	53	106	74-149
Tetrahydrofuran	50	NA	46	93	46-154
1,1,1-Trichloroethane	50	NA	52	104	76-145
1,1-Dichloropropene	50	NA	51	102	87-130
2-Butanone	50	NA	51	101	30-150
Benzene	50	NA	46	93	90-116
Cyclohexane	50	NA	47	94	70-130
Ethyl Methacrylate	50	NA	55	109	60-138
Tertiary-amyl methyl ether	50	NA	52	104	82-118
1,2-Dichloroethane	50	NA	48	96	71-149
Trichloroethene	50	NA	48	97	83-132

KATAHDIN ANALYTICAL SERVICES
LAB CONTROL SAMPLE

Client:
Project: Old Town
PO No:
Sample Date:
Received Date:
Extraction Date:
Analysis Date: 05/31/06
Report Date: 06/07/2006
Matrix: WATER

Lab ID: WG28646-1
Client ID: WG28646-LCS
SDG: WW2637
Extracted by:
Extraction Method: SW846 5030
Analyst: SKT
Analysis Method: SW846 8260B
Lab Prep Batch: WG28646
Units: ug/l

COMPOUND	LCS SPIKE	SAMPLE CONC.	LCS CONC.		%REC.	QC. LIMITS
Dibromomethane	50	NA	46		92	76-132
1,2-Dichloropropane	50	NA	47		95	95-112
Bromodichloromethane	50	NA	46		92	76-132
cis-1,3-dichloropropene	50	NA	50		100	85-127
1,4-Dioxane	1000	NA	1000		100	0-125
2-Chloroethylvinylether	50	NA	77	*	154	20-137
Toluene	50	NA	49		97	85-122
4-methyl-2-pentanone	50	NA	55		110	63-154
Tetrachloroethene	50	NA	52		105	48-155
trans-1,3-Dichloropropene	50	NA	54		108	85-144
1,1,2-Trichloroethane	50	NA	48		97	78-124
Dibromochloromethane	50	NA	49		98	63-128
1,3-Dichloropropane	50	NA	48		96	65-127
1,2-Dibromoethane	50	NA	50		101	75-129
2-Hexanone	50	NA	54		108	47-160
Chlorobenzene	50	NA	50		99	68-128
Ethylbenzene	50	NA	49		98	76-124
1,1,1,2-Tetrachloroethane	50	NA	51		102	69-133
Xylenes (total)	150	NA	157		105	67-129
m+p-Xylenes	100	NA	105		105	68-127
o-Xylene	50	NA	52		104	64-133
Styrene	50	NA	52		105	66-130
Bromoform	50	NA	54		108	59-135
Isopropylbenzene	50	NA	58		117	95-140
cis-1,4-Dichloro-2-Butene	50	NA	44		89	56-155
trans-1,4-Dichloro-2-Butene	50	NA	46		91	66-151
Bromobenzene	50	NA	48		96	86-122
N-Propylbenzene	50	NA	53		106	92-124
1,1,2,2-Tetrachloroethane	50	NA	51		102	75-128
1,3,5-Trimethylbenzene	50	NA	54		108	86-127
2-Chlorotoluene	50	NA	51		101	84-129
1,2,3-Trichloropropane	50	NA	51		102	57-145
4-Chlorotoluene	50	NA	51		102	83-129
tert-Butylbenzene	50	NA	55		110	86-131
Pentachloroethane	50	NA	51		102	22-197
1,2,4-Trimethylbenzene	50	NA	53		106	85-124
P-Isopropyltoluene	50	NA	58		115	84-131
1,3-Dichlorobenzene	50	NA	51		103	81-121
1,4-Dichlorobenzene	50	NA	57		115	81-120
N-Butylbenzene	50	NA	57		114	67-138
sec-Butylbenzene	50	NA	57		114	82-128
1,2-Dichlorobenzene	50	NA	53		105	74-124
1,2-Dibromo-3-Chloropropane	50	NA	52		103	55-151
1,3,5-Trichlorobenzene	50	NA	50		101	77-122
Hexachlorobutadiene	50	NA	51		102	75-119

KATAHDIN ANALYTICAL SERVICES
LAB CONTROL SAMPLE

Client:
Project: Old Town
PO No:
Sample Date:
Received Date:
Extraction Date:
Analysis Date: 05/31/06
Report Date: 06/07/2006
Matrix: WATER

Lab ID: WG28646-1
Client ID: WG28646-LCS
SDG: WW2637
Extracted by:
Extraction Method: SW846 5030
Analyst: SKT
Analysis Method: SW846 8260B
Lab Prep Batch: WG28646
Units: ug/l

COMPOUND	LCS SPIKE	SAMPLE CONC.	LCS CONC.	%REC.	QC. LIMITS
1,2,4-Trichlorobenzene	50	NA	54	109	64-125
1,2,3-Trimethylbenzene	50	NA	52	104	84-122
Naphthalene	50	NA	45	90	40-137
1,2,3-Trichlorobenzene	50	NA	48	96	38-140
Methyl Acetate	50	NA	52	104	61-168
Methylcyclohexane	50	NA	54	107	81-130
1-Chlorohexane	50	NA	53	106	80-120
Total Alkylbenzenes	400	NA	445	111	60-140

KATAHDIN ANALYTICAL SERVICES
MATRIX SPIKE/MATRIX SPIKE DUPLICATE RECOVERY

Client: Woodard & Curran
Project: Old Town
PO No:
Sample Date: 05/30/06
Received Date: 05/31/06
Extraction Date:
Analysis Date: 05/31/06
Report Date: 06/07/2006
Matrix: WATER

Lab ID: WG28646-3 & WG28646-4
Client ID: SB-25MS & SB-25MSD
SDG: WW2637
Extracted by:
Extraction Method: SW846 5030
Analyst: SKT
Analysis Method: SW846 8260B
Lab Prep Batch: WG28646
Units: ug/l

COMPOUND	MS SPIKE	MSD SPIKE	SAMPLE CONC.	MS CONC.	MSD CONC.	MS %REC.	MSD %REC.	%RPD	%RPD LIMIT	QC. LIMITS
Dichlorodifluoromethane	50	50	0.00	39	37	77	74	4	20	0-200
Chloromethane	50	50	0.00	48	46	95	93	2	20	40-125
Vinyl chloride	50	50	0.00	51	53	102	107	5	20	27-143
Bromomethane	50	50	0.00	33	48	66	96 *	38	20	52-141
Chloroethane	50	50	0.00	57	58	114	115	1	20	64-124
Trichlorofluoromethane	50	50	0.00	79	78	* 158	* 157	0.9	20	63-153
1,1-Dichloroethene	50	50	0.00	52	56	105	113	8	20	85-129
Methylene Chloride	50	50	0.00	51	51	103	103	0.0	20	86-124
trans-1,2-Dichloroethene	50	50	0.00	49	50	98	100	2	20	80-128
1,1-Dichloroethane	50	50	0.00	60	59	120	118	2	20	90-127
cis-1,2-Dichloroethene	50	50	0.00	46	48	93	96	4	20	84-121
1,2-Dichloroethylene (total)	100	100	0.00	95	98	95	98	3	20	83-123
2,2-Dichloropropane	50	50	0.00	48	47	95	93	2	20	68-149
Chloroform	50	50	0.00	57	57	114	113	0.7	20	87-127
Bromochloromethane	50	50	0.00	53	53	105	106	0.4	20	87-120
1,1,1-Trichloroethane	50	50	0.00	61	60	122	120	1	20	76-145
1,2-Dichloroethane	50	50	0.00	55	54	109	109	0.5	20	71-149
1,1-Dichloropropene	50	50	0.00	59	60	119	120	1	20	87-130
Carbon Tetrachloride	50	50	0.00	60	60	121	120	0.8	20	74-149
Benzene	50	50	0.00	54	54	108	108	0.6	20	90-116
1,2-Dichloropropane	50	50	0.00	58	57	* 117	* 114	3	20	95-112
Trichloroethene	50	50	0.00	53	53	106	107	0.9	20	83-132
Dibromomethane	50	50	0.00	53	51	106	102	3	20	76-132
Bromodichloromethane	50	50	0.00	52	51	103	102	1.0	20	76-132
cis-1,3-dichloropropene	50	50	0.00	56	56	111	111	0.0	20	85-127
Toluene	50	50	0.00	52	52	103	105	1	20	85-122
trans-1,3-Dichloropropene	50	50	0.00	61	60	122	120	1	20	85-144
1,1,2-Trichloroethane	50	50	0.00	54	53	108	106	2	20	78-124
1,3-Dichloropropane	50	50	0.00	52	52	104	104	0.4	20	65-127
Dibromochloromethane	50	50	0.00	48	49	95	97	2	20	63-128
Tetrachloroethene	50	50	0.00	51	52	101	105	3	20	48-155
1,2-Dibromoethane	50	50	0.00	54	54	108	109	1	20	75-129
Chlorobenzene	50	50	0.00	49	50	98	101	3	20	68-128
1,1,1,2-Tetrachloroethane	50	50	0.00	51	52	101	105	3	20	69-133
Ethylbenzene	50	50	0.00	50	51	100	102	2	20	76-124
Bromoform	50	50	0.00	52	54	104	107	3	20	59-135
Styrene	50	50	0.00	51	53	101	106	4	20	66-130
1,1,2,2-Tetrachloroethane	50	50	0.00	56	56	112	111	0.4	20	75-128
1,2,3-Trichloropropane	50	50	0.00	57	55	114	110	3	20	57-145
Isopropylbenzene	50	50	0.00	58	60	117	121	4	20	95-140
Bromobenzene	50	50	0.00	47	50	94	100	6	20	86-122
2-Chlorotoluene	50	50	0.00	51	52	103	104	1	20	84-129
N-Propylbenzene	50	50	0.00	55	56	109	112	2	20	92-124
4-Chlorotoluene	50	50	0.00	52	53	103	106	3	20	83-129
1,3,5-Trimethylbenzene	50	50	0.00	55	56	109	111	2	20	86-127

FORM III VOA-1
T1148.D & T1149.D

KATAHDIN ANALYTICAL SERVICES
MATRIX SPIKE/MATRIX SPIKE DUPLICATE RECOVERY

Client: Woodard & Curran
Project: Old Town
PO No:
Sample Date: 05/30/06
Received Date: 05/31/06
Extraction Date:
Analysis Date: 05/31/06
Report Date: 06/07/2006
Matrix: WATER

Lab ID: WG28646-3 & WG28646-4
Client ID: SB-25MS & SB-25MSD
SDG: WW2637
Extracted by:
Extraction Method: SW846 5030
Analyst: SKT
Analysis Method: SW846 8260B
Lab Prep Batch: WG28646
Units: ug/l

COMPOUND	MS SPIKE	MSD SPIKE	SAMPLE CONC.	MS CONC.	MSD CONC.	MS %REC.	MSD %REC.	%RPD	%RPD LIMIT	QC. LIMITS
tert-Butylbenzene	50	50	0.00	56	58	112	115	3	20	86-131
1,2,4-Trichlorobenzene	50	50	0.00	55	56	110	111	0.7	20	64-125
sec-Butylbenzene	50	50	0.00	58	59	116	118	2	20	82-128
1,3-Dichlorobenzene	50	50	0.00	52	52	103	105	2	20	81-121
P-Isopropyltoluene	50	50	0.00	59	60	117	119	2	20	84-131
1,4-Dichlorobenzene	50	50	0.00	58	60	117	120	3	20	81-120
1,2-Dichlorobenzene	50	50	0.00	53	54	105	107	2	20	74-124
N-Butylbenzene	50	50	0.00	59	58	117	117	0.3	20	67-138
1,2-Dibromo-3-Chloropropane	50	50	0.00	57	58	114	115	0.5	20	55-151
1,2,4-Trimethylbenzene	50	50	0.00	52	54	105	108	3	20	85-124
Naphthalene	50	50	0.00	48	47	96	94	2	20	40-137
Hexachlorobutadiene	50	50	0.00	50	51	101	103	2	20	75-119
1,2,3-Trichlorobenzene	50	50	0.00	51	50	102	100	2	20	38-140
Methyl tert-butyl ether	100	100	0.00	116	121	116	121	4	20	79-132
Acetone	50	50	0.00	74	71	149	142	4	20	41-180
2-Butanone	50	50	0.00	64	60	127	119	6	20	30-150
4-methyl-2-pentanone	50	50	0.00	70	67	140	134	4	20	63-154
2-Hexanone	50	50	0.00	63	61	127	122	3	20	47-160
m+p-Xylenes	100	100	0.00	104	107	104	107	3	20	68-127
o-Xylene	50	50	0.00	51	53	101	105	4	20	64-133
Xylenes (total)	150	150	0.00	155	160	103	107	3	20	67-129
1,3,5-Trichlorobenzene	50	50	0.00	47	53	94	106	12	20	77-122
Vinyl Acetate	50	50	0.00	53	53	105	106	0.6	20	47-122
Carbon Disulfide	50	50	0.00	56	60	113	120	6	20	80-134
Diethyl Ether	50	50	0.00	48	58	95	116	19	20	63-146
Tetrahydrofuran	50	50	0.00	56	54	112	108	4	20	46-154

FORM III VOA-1
T1148.D & T1149.D

FORM 4

VOLATILE METHOD BLANK SUMMARY

CLIENT SAMPLE ID

| WG28636-BLANK |

Lab Name: KATAHDIN ANALYTICAL SERVICES Lab Code: KAS

Project: OLD TOWN SDG No.: WW2637

Lab File ID: Z0278 Lab Sample ID: WG28636-2

Date Analyzed: 05/31/06 Time Analyzed: 1002

GC Column: RTX-VMS ID: 0.18 (mm) Heated Purge: (Y/N) Y

Instrument ID: GCMS-Z

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES, MS and MSD:

	CLIENT SAMPLE ID	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED	TIME ANALYZED
01	WG28636-LCS	WG28636-1	Z0276	05/31/06	0842
02	SS01	WW2637-6	Z0280	05/31/06	1157
03	SS04	WW2637-10	Z0282	05/31/06	1310
04	SS05	WW2637-11	Z0283	05/31/06	1347
05	SS81	WW2637-14	Z0284	05/31/06	1423
06	SS03	WW2637-8RA	Z0285	05/31/06	1501
07	SS05MS	WG28636-3	Z0288	05/31/06	1651
08	SS05MSD	WG28636-4	Z0289	05/31/06	1727
09					
10					
11					
12					
13					
14					
15					
16					
17					
18					
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20					
21					
22					
23					
24					
25					
26					
27					
28					
29					
30					

COMMENTS:

KATAHDIN ANALYTICAL SERVICES
Report of Analytical Results

Client:
Project: Old Town
PO No:
Sample Date:
Received Date:
Extraction Date:
Analysis Date: 31-MAY-2006 10:02
Report Date: 06/07/2006
Matrix: SOIL
% Solids: 100

Lab ID: WG28636-2
Client ID: WG28636-Blank
SDG: WW2637
Extracted by:
Extraction Method: SW846 5035
Analyst: ALH
Analysis Method: SW846 8260B
Lab Prep Batch: WG28636
Units: ug/Kg

Compound	Flags	Results	DF	PQL	Adj.PQL
Dichlorodifluoromethane	U	5	1.0	5	5
Chloromethane	U	5	1.0	5	5
Vinyl chloride	U	5	1.0	5	5
Bromomethane	U	5	1.0	5	5
Chloroethane	U	5	1.0	5	5
Trichlorofluoromethane	U	5	1.0	5	5
1,1-Dichloroethene	U	5	1.0	5	5
Methylene Chloride	U	25	1.0	25	25
trans-1,2-Dichloroethene	U	5	1.0	5	5
1,1-Dichloroethane	U	5	1.0	5	5
cis-1,2-Dichloroethene	U	5	1.0	5	5
1,2-Dichloroethylene (total)	U	10	1.0	10	10
2,2-Dichloropropane	U	5	1.0	5	5
Chloroform	U	5	1.0	5	5
Bromochloromethane	U	5	1.0	5	5
1,1,1-Trichloroethane	U	5	1.0	5	5
1,2-Dichloroethane	U	5	1.0	5	5
1,1-Dichloropropene	U	5	1.0	5	5
Carbon Tetrachloride	U	5	1.0	5	5
Benzene	U	5	1.0	5	5
1,2-Dichloropropane	U	5	1.0	5	5
Trichloroethene	U	5	1.0	5	5
Dibromomethane	U	5	1.0	5	5
Bromodichloromethane	U	5	1.0	5	5
cis-1,3-dichloropropene	U	5	1.0	5	5
Toluene	U	5	1.0	5	5
trans-1,3-Dichloropropene	U	5	1.0	5	5
1,1,2-Trichloroethane	U	5	1.0	5	5
1,3-Dichloropropane	U	5	1.0	5	5
Dibromochloromethane	U	5	1.0	5	5
Tetrachloroethene	U	5	1.0	5	5
1,2-Dibromoethane	U	5	1.0	5	5
Chlorobenzene	U	5	1.0	5	5
1,1,1,2-Tetrachloroethane	U	5	1.0	5	5
Ethylbenzene	U	5	1.0	5	5
Bromoform	U	5	1.0	5	5
Styrene	U	5	1.0	5	5
1,1,2,2-Tetrachloroethane	U	5	1.0	5	5
1,2,3-Trichloropropane	U	5	1.0	5	5
Isopropylbenzene	U	5	1.0	5	5
Bromobenzene	U	5	1.0	5	5
2-Chlorotoluene	U	5	1.0	5	5
N-Propylbenzene	U	5	1.0	5	5

KATAHDIN ANALYTICAL SERVICES
Report of Analytical Results

Client:
Project: Old Town
PO No:
Sample Date:
Received Date:
Extraction Date:
Analysis Date: 31-MAY-2006 10:02
Report Date: 06/07/2006
Matrix: SOIL
% Solids: 100

Lab ID: WG28636-2
Client ID: WG28636-Blank
SDG: WW2637
Extracted by:
Extraction Method: SW846 5035
Analyst: ALH
Analysis Method: SW846 8260B
Lab Prep Batch: WG28636
Units: ug/Kg

Compound	Flags	Results	DF	PQL	Adj.PQL
4-Chlorotoluene	U	5	1.0	5	5
1,3,5-Trimethylbenzene	U	5	1.0	5	5
tert-Butylbenzene	U	5	1.0	5	5
1,2,4-Trichlorobenzene	U	5	1.0	5	5
sec-Butylbenzene	U	5	1.0	5	5
1,3-Dichlorobenzene	U	5	1.0	5	5
P-Isopropyltoluene	U	5	1.0	5	5
1,4-Dichlorobenzene	U	5	1.0	5	5
1,2-Dichlorobenzene	U	5	1.0	5	5
N-Butylbenzene	U	5	1.0	5	5
1,2-Dibromo-3-Chloropropane	U	5	1.0	5	5
1,2,4-Trimethylbenzene	U	5	1.0	5	5
Naphthalene		5	1.0	5	5
Hexachlorobutadiene	U	5	1.0	5	5
1,2,3-Trichlorobenzene	U	5	1.0	5	5
Methyl tert-butyl ether	U	5	1.0	5	5
Acetone	U	25	1.0	25	25
2-Butanone	U	25	1.0	25	25
4-methyl-2-pentanone	U	25	1.0	25	25
2-Hexanone	U	25	1.0	25	25
m+p-Xylenes	U	10	1.0	10	10
o-Xylene	U	5	1.0	5	5
Xylenes (total)	U	15	1.0	15	15
1,3,5-Trichlorobenzene	U	5	1.0	5	5
Vinyl Acetate	U	5	1.0	5	5
Carbon Disulfide	U	5	1.0	5	5
Diethyl Ether	U	5	1.0	5	5
Tetrahydrofuran	U	25	1.0	25	25
Dibromofluoromethane		81%			
1,2-Dichloroethane-D4		77%			
Toluene-D8		82%			
P-Bromofluorobenzene		74%			

KATAHDIN ANALYTICAL SERVICES
LAB CONTROL SAMPLE

Client:
Project: Old Town
PO No:
Sample Date:
Received Date:
Extraction Date:
Analysis Date: 05/31/06
Report Date: 06/07/2006
Matrix: SOIL

Lab ID: WG28636-1
Client ID: WG28636-LCS
SDG: WW2637
Extracted by:
Extraction Method: SW846 5035
Analyst: ALH
Analysis Method: SW846 8260B
Lab Prep Batch: WG28636
Units: ug/Kg

COMPOUND	LCS SPIKE	SAMPLE CONC.	LCS CONC.		%REC.	QC. LIMITS
Dichlorodifluoromethane	50	NA	25		50	0-200
Chloromethane	50	NA	34		69	50-130
Vinyl chloride	50	NA	37		74	16-149
Bromomethane	50	NA	62		124	29-151
Chloroethane	50	NA	44		87	50-127
Trichlorofluoromethane	50	NA	52		104	56-141
Diethyl Ether	50	NA	53		107	53-139
Tertiary-butyl alcohol	250	NA	134	*	54	60-146
1,1-Dichloroethene	50	NA	42		84	68-134
Carbon Disulfide	50	NA	48		96	54-154
Freon-113	50	NA	57		114	73-131
Iodomethane	50	NA	49		98	8-175
Acrolein	250	NA	238		95	46-154
Methylene Chloride	50	NA	46		92	65-139
Acetone	50	NA	52		105	20-160
Isobutyl Alcohol	1000	NA	720		72	50-145
trans-1,2-Dichloroethene	50	NA	47		93	67-139
Allyl Chloride	50	NA	60		121	51-180
Methyl tert-butyl ether	100	NA	91		91	83-128
Acetonitrile	500	NA	425		85	31-157
Di-isopropyl ether	50	NA	48		97	69-136
Chloroprene	50	NA	50		101	83-121
Methacrylonitrile	500	NA	494		99	66-137
Propionitrile	500	NA	451		90	38-157
1,1-Dichloroethane	50	NA	48		96	74-129
Acrylonitrile	250	NA	230		92	42-155
Ethyl tertiary-butyl ether	50	NA	50		101	77-129
Vinyl Acetate	50	NA	42		84	64-102
cis-1,2-Dichloroethene	50	NA	44		87	72-130
1,2-Dichloroethylene (total)	100	NA	90		90	70-134
Methyl Methacrylate	50	NA	47		95	80-130
2,2-Dichloropropane	50	NA	59		117	76-132
Bromochloromethane	50	NA	46		92	81-127
Chloroform	50	NA	47		94	80-125
Carbon Tetrachloride	50	NA	52		104	69-148
Tetrahydrofuran	50	NA	44		87	64-147
1,1,1-Trichloroethane	50	NA	49		98	83-130
1,1-Dichloropropene	50	NA	49		97	86-127
2-Butanone	50	NA	67		135	30-160
Benzene	50	NA	50		100	81-121
Cyclohexane	50	NA	42		83	64-138
Ethyl Methacrylate	50	NA	48		96	76-129
Tertiary-amyl methyl ether	50	NA	49		99	76-126
1,2-Dichloroethane	50	NA	49		98	69-148
Trichloroethene	50	NA	52		104	86-121

KATAHDIN ANALYTICAL SERVICES
LAB CONTROL SAMPLE

Client:
Project: Old Town
PO No:
Sample Date:
Received Date:
Extraction Date:
Analysis Date: 05/31/06
Report Date: 06/07/2006
Matrix: SOIL

Lab ID: WG28636-1
Client ID: WG28636-LCS
SDG: WW2637
Extracted by:
Extraction Method: SW846 5035
Analyst: ALH
Analysis Method: SW846 8260B
Lab Prep Batch: WG28636
Units: ug/Kg

COMPOUND	LCS SPIKE	SAMPLE CONC.	LCS CONC.	%REC.	QC. LIMITS
Dibromomethane	50	NA	50	100	75-131
1,2-Dichloropropane	50	NA	49	98	78-125
Bromodichloromethane	50	NA	48	95	78-123
cis-1,3-dichloropropene	50	NA	53	106	91-121
1,4-Dioxane	1000	NA	819	82	61-151
2-Chloroethylvinylether	50	NA	38	76	30-135
Toluene	50	NA	54	107	86-120
4-methyl-2-pentanone	50	NA	48	96	79-130
Tetrachloroethene	50	NA	54	108	74-136
trans-1,3-Dichloropropene	50	NA	59	117	88-134
1,1,2-Trichloroethane	50	NA	50	100	85-116
Dibromochloromethane	50	NA	51	101	81-120
1,3-Dichloropropane	50	NA	50	101	90-112
1,2-Dibromoethane	50	NA	51	102	89-119
2-Hexanone	50	NA	51	102	84-132
Chlorobenzene	50	NA	54	108	86-114
Ethylbenzene	50	NA	55	109	88-114
1,1,1,2-Tetrachloroethane	50	NA	54	109	83-126
Xylenes (total)	150	NA	162	108	84-118
m+p-Xylenes	100	NA	108	108	84-119
o-Xylene	50	NA	54	108	82-120
Styrene	50	NA	54	107	85-115
Bromoform	50	NA	58	116	84-124
Isopropylbenzene	50	NA	57	114	92-126
cis-1,4-Dichloro-2-Butene	50	NA	47	94	75-128
trans-1,4-Dichloro-2-Butene	50	NA	48	97	73-125
Bromobenzene	50	NA	52	104	84-116
N-Propylbenzene	50	NA	53	107	81-121
1,1,2,2-Tetrachloroethane	50	NA	48	97	74-128
1,3,5-Trimethylbenzene	50	NA	52	105	87-114
2-Chlorotoluene	50	NA	50	100	83-117
1,2,3-Trichloropropane	50	NA	54	107	38-138
4-Chlorotoluene	50	NA	52	103	84-117
tert-Butylbenzene	50	NA	54	108	72-128
Pentachloroethane	50	NA	52	104	75-119
1,2,4-Trimethylbenzene	50	NA	52	104	88-113
P-Isopropyltoluene	50	NA	56	112	85-123
1,3-Dichlorobenzene	50	NA	56	112	84-118
1,4-Dichlorobenzene	50	NA	56	113	87-117
N-Butylbenzene	50	NA	53	106	83-116
sec-Butylbenzene	50	NA	53	106	87-115
1,2-Dichlorobenzene	50	NA	56	111	86-114
1,2-Dibromo-3-Chloropropane	50	NA	48	97	83-117
1,3,5-Trichlorobenzene	50	NA	55	111	67-127
Hexachlorobutadiene	50	NA	58	116	77-122

KATAHDIN ANALYTICAL SERVICES
LAB CONTROL SAMPLE

Client:
Project: Old Town
PO No:
Sample Date:
Received Date:
Extraction Date:
Analysis Date: 05/31/06
Report Date: 06/07/2006
Matrix: SOIL

Lab ID: WG28636-1
Client ID: WG28636-LCS
SDG: WW2637
Extracted by:
Extraction Method: SW846 5035
Analyst: ALH
Analysis Method: SW846 8260B
Lab Prep Batch: WG28636
Units: ug/Kg

COMPOUND	LCS SPIKE	SAMPLE CONC.	LCS CONC.	%REC.	QC. LIMITS
1,2,4-Trichlorobenzene	50	NA	52	103	74-124
1,2,3-Trimethylbenzene	50	NA	53	106	84-116
Naphthalene	50	NA	35	70	29-166
1,2,3-Trichlorobenzene	50	NA	43	86	45-155
Methyl Acetate	50	NA	58	116	36-185
Methylcyclohexane	50	NA	54	107	68-131

KATAHDIN ANALYTICAL SERVICES
MATRIX SPIKE/MATRIX SPIKE DUPLICATE RECOVERY

Client: Woodard & Curran
Project: Old Town
PO No:
Sample Date: 05/30/06
Received Date: 05/31/06
Extraction Date:
Analysis Date: 05/31/06
Report Date: 06/07/2006
Matrix: SOIL

Lab ID: WG28636-3 & WG28636-4
Client ID: SS05MS & SS05MSD
SDG: WW2637
Extracted by:
Extraction Method: SW846 5035
Analyst: ALH
Analysis Method: SW846 8260B
Lab Prep Batch: WG28636
Units: ug/Kg

COMPOUND	MS SPIKE	MSD SPIKE	SAMPLE CONC.	MS CONC.	MSD CONC.		MS %REC.		MSD %REC.		%RPD	%RPD LIMIT	QC. LIMITS
Dichlorodifluoromethane	31	32	0.00	10	15		33		48	*	39	30	0-200
Chloromethane	31	32	0.00	8.9	14	*	29	*	44	*	44	30	50-130
Vinyl chloride	31	32	0.00	13	20		43		64	*	43	30	16-149
Bromomethane	31	32	0.00	15	21		49		66	*	32	30	29-151
Chloroethane	31	32	0.00	17	23		54		73	*	32	30	50-127
Trichlorofluoromethane	31	32	0.00	20	31		66		98	*	43	30	56-141
1,1-Dichloroethene	31	32	0.00	15	22	*	49		70	*	39	30	68-134
Methylene Chloride	31	32	0.00	15	18	*	49	*	55		15	30	65-139
trans-1,2-Dichloroethene	31	32	0.00	14	20	*	44	*	63	*	38	30	67-139
1,1-Dichloroethane	31	32	0.00	17	21	*	56	*	67		21	30	74-129
cis-1,2-Dichloroethene	31	32	0.00	14	17	*	46	*	54		20	30	72-130
1,2-Dichloroethylene (total)	62	63	0.00	28	37	*	45	*	59		30	30	70-134
2,2-Dichloropropane	31	32	0.00	10	13	*	34	*	40		20	30	76-132
Chloroform	31	32	0.00	16	19	*	54	*	61		16	30	80-125
Bromochloromethane	31	32	0.00	16	17	*	53	*	53		2	30	81-127
1,1,1-Trichloroethane	31	32	0.00	16	24	*	53	*	75	*	37	30	83-130
1,2-Dichloroethane	31	32	0.00	17	16	*	56	*	51		7	30	69-148
1,1-Dichloropropene	31	32	0.00	14	22	*	46	*	69	*	42	30	86-127
Carbon Tetrachloride	31	32	0.00	16	24	*	52		76	*	40	30	69-148
Benzene	31	32	0.00	15	20	*	50	*	62		23	30	81-121
1,2-Dichloropropane	31	32	0.00	15	17	*	50	*	53		10	30	78-125
Trichloroethene	31	32	0.00	14	20	*	45	*	62	*	33	30	86-121
Dibromomethane	31	32	0.00	15	15	*	49	*	47		1	30	75-131
Bromodichloromethane	31	32	0.00	14	14	*	44	*	44		4	30	78-123
cis-1,3-dichloropropene	31	32	0.00	10	11	*	34	*	33		2	30	91-121
Toluene	31	32	0.00	13	17	*	41	*	54		29	30	86-120
trans-1,3-Dichloropropene	31	32	0.00	11	10	*	34	*	32		5	30	88-134
1,1,2-Trichloroethane	31	32	0.00	14	14	*	47	*	44		4	30	85-116
1,3-Dichloropropane	31	32	0.00	15	15	*	48	*	46		2	30	90-112
Dibromochloromethane	31	32	0.00	12	12	*	38	*	37		2	30	81-120
Tetrachloroethene	31	32	0.00	13	20	*	43	*	65	*	43	30	74-136
1,2-Dibromoethane	31	32	0.00	13	12	*	41	*	38		6	30	89-119
Chlorobenzene	31	32	0.00	11	13	*	36	*	42		19	30	86-114
1,1,1,2-Tetrachloroethane	31	32	0.00	13	14	*	42	*	44		10	30	83-126
Ethylbenzene	31	32	0.00	11	16	*	36	*	52	*	40	30	88-114
Bromoform	31	32	0.00	9.7	9.5	*	32	*	30		2	30	84-124
Styrene	31	32	0.00	7.4	8.3	*	24	*	26	*	11	3	85-115
1,1,2,2-Tetrachloroethane	31	32	0.00	13	12	*	42	*	37		8	30	74-128
1,2,3-Trichloropropane	31	32	0.00	6.6	6.0	*	21	*	19		10	30	38-138
Isopropylbenzene	31	32	0.00	12	20	*	41	*	64	*	47	30	92-126
Bromobenzene	31	32	0.00	9.8	11	*	32	*	34		11	30	84-116
2-Chlorotoluene	31	32	0.00	9.0	13	*	29	*	42	*	39	30	83-117
N-Propylbenzene	31	32	0.00	10	16	*	33	*	50	*	45	30	81-121
4-Chlorotoluene	31	32	0.00	9.2	11	*	30	*	35		20	30	84-117
1,3,5-Trimethylbenzene	31	32	0.00	9.7	14	*	32	*	43	*	33	30	87-114

FORM III VOA-2
Z0288.D & Z0289.D

KATAHDIN ANALYTICAL SERVICES
MATRIX SPIKE/MATRIX SPIKE DUPLICATE RECOVERY

Client: Woodard & Curran
Project: Old Town
PO NO:
Sample Date: 05/30/06
Received Date: 05/31/06
Extraction Date:
Analysis Date: 05/31/06
Report Date: 06/07/2006
Matrix: SOIL

Lab ID: WG28636-3 & WG28636-4
Client ID: SS05MS & SS05MSD
SDG: WW2637
Extracted by:
Extraction Method: SW846 5035
Analyst: ALH
Analysis Method: SW846 8260B
Lab Prep Batch: WG28636
Units: ug/Kg

COMPOUND	MS SPIKE	MSD SPIKE	SAMPLE CONC.	MS CONC.	MSD CONC.		MS %REC.		MSD %REC.		%RPD	%RPD LIMIT	QC. LIMITS
tert-Butylbenzene	31	32	0.00	11	16	*	34	*	50	*	39	30	72-128
1,2,4-Trichlorobenzene	31	32	0.00	3.5	2.0	*	11	*	6	*	54	30	74-124
sec-Butylbenzene	31	32	0.00	10	16	*	33	*	51	*	46	30	87-115
1,3-Dichlorobenzene	31	32	0.00	8.4	9.0	*	27	*	28		7	30	84-118
P-Isopropyltoluene	31	32	0.00	8.9	13	*	29	*	41	*	38	30	85-123
1,4-Dichlorobenzene	31	32	0.00	8.2	8.7	*	27	*	27		6	30	87-117
1,2-Dichlorobenzene	31	32	0.00	7.5	7.4	*	24	*	23		1	30	86-114
N-Butylbenzene	31	32	0.00	7.5	9.1	*	24	*	29		19	30	83-116
1,2-Dibromo-3-Chloropropane	31	32	0.00	6.0	4.9	*	19	*	15		20	30	83-117
1,2,4-Trimethylbenzene	31	32	0.00	8.7	12	*	28	*	36		28	30	88-113
Naphthalene	31	32	0.00	4.6	4.0	*	15	*	13		14	30	29-166
Hexachlorobutadiene	31	32	0.00	7.1	7.3	*	23	*	23		3	30	77-122
1,2,3-Trichlorobenzene	31	32	0.00	4.9	3.7	*	16	*	12		28	30	45-155
Methyl tert-butyl ether	62	63	0.00	34	31	*	55	*	49		8	30	83-128
Acetone	31	32	71	472	435	*	1303	*	1149		8	30	20-160
2-Butanone	31	32	0.00	27	25		88		79		8	30	30-160
4-methyl-2-pentanone	31	32	0.00	12	8.5	*	40	*	27	*	36	30	79-130
2-Hexanone	31	32	0.00	10	6.7	*	33	*	21	*	41	30	84-132
m+p-Xylenes	62	63	0.00	21	29	*	34	*	46	*	33	30	84-119
o-Xylene	31	32	0.00	9.9	13	*	32	*	41		27	30	82-120
Xylenes (total)	92	95	0.00	31	42	*	33	*	44	*	31	30	84-118
1,3,5-Trichlorobenzene	31	32	0.00	4.2	3.3	*	14	*	10		24	30	67-127
Vinyl Acetate	31	32	0.00	5.6	5.7	*	18	*	18		2	30	64-102
Carbon Disulfide	31	32	0.00	8.5	15	*	28	*	46	*	53	30	54-154
Diethyl Ether	31	32	0.00	20	22		65		68		8	30	53-139
Tetrahydrofuran	31	32	0.00	16	14	*	51	*	46		8	30	64-147

FORM III VOA-2
Z0288.D & Z0289.D

FORM 4

SEMIVOLATILE METHOD BLANK SUMMARY

CLIENT SAMPLE ID

| WG28621-BLANK |

Lab Name: KATAHDIN ANALYTICAL SERVICES Lab Code: KAS

Project: OLD TOWN SDG No.: WW2637

Lab File ID: R1171 Lab Sample ID: WG28621-1

Instrument ID: GCMS-R Date Extracted: 05/31/06

Matrix: (soil/water) SOIL Date Analyzed: 06/01/06

Level:(low/med) LOW Time Analyzed: 1126

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES, MS and MSD:

	CLIENT SAMPLE ID	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED	TIME ANALYZED
01	SS01	WW2637-6	R1161	05/31/06	1713
02	SS03	WW2637-8	R1162	05/31/06	1759
03	SS80	WW2637-9	R1163	05/31/06	1844
04	SS04	WW2637-10	R1164	05/31/06	1928
05	WG28621-LCS	WG28621-2	R1172	06/01/06	1212
06	WG28621-LCSD	WG28621-3	R1173	06/01/06	1258
07	SS05	WW2637-11RA	R1182	06/02/06	0944
08	SS02	WW2637-7DL2	R1183	06/02/06	1031
09					
10					
11					
12					
13					
14					
15					
16					
17					
18					
19					
20					
21					
22					
23					
24					
25					
26					
27					
28					
29					
30					

COMMENTS:

FORM IV SV

KATAHDIN ANALYTICAL SERVICES
Report of Analytical Results

Client:
Project: Old Town
PO No:
Sample Date:
Received Date:
Extraction Date: 05/31/06
Analysis Date: 01-JUN-2006 11:26
Report Date: 06/06/2006
Matrix: SOIL
% Solids: 100

Lab ID: WG28621-1
Client ID: WG28621-Blank
SDG: WW2637
Extracted by: KM
Extraction Method: SW846 3550
Analyst: LRS
Analysis Method: SW846 8270C
Lab Prep Batch: WG28621
Units: ug/Kg

Compound	Flags	Results	DF	PQL	Adj.PQL
Phenol	U	330	1.0	330	330
Bis(2-Chloroethyl)ether	U	330	1.0	330	330
2-Chlorophenol	U	330	1.0	330	330
1,3-Dichlorobenzene	U	330	1.0	330	330
1,4-Dichlorobenzene	U	330	1.0	330	330
1,2-Dichlorobenzene	U	330	1.0	330	330
2-Methylphenol	U	330	1.0	330	330
2,2'-Oxybis(1-Chloropropane)	U	330	1.0	330	330
N-Nitroso-di-n-propylamine	U	330	1.0	330	330
3&4-Methylphenol	U	330	1.0	330	330
Hexachloroethane	U	330	1.0	330	330
Nitrobenzene	U	330	1.0	330	330
Isophorone	U	330	1.0	330	330
2-Nitrophenol	U	330	1.0	330	330
2,4-Dimethylphenol	U	330	1.0	330	330
Bis(2-Chloroethoxy)methane	U	330	1.0	330	330
2,4-Dichlorophenol	U	330	1.0	330	330
1,2,4-Trichlorobenzene	U	330	1.0	330	330
Naphthalene	U	330	1.0	330	330
4-Chloroaniline	U	330	1.0	330	330
Hexachlorobutadiene	U	330	1.0	330	330
4-Chloro-3-Methylphenol	U	330	1.0	330	330
2-Methylnaphthalene	U	330	1.0	330	330
Hexachlorocyclopentadiene	U	330	1.0	330	330
2,4,6-Trichlorophenol	U	330	1.0	330	330
2,4,5-Trichlorophenol	U	820	1.0	820	820
2-Chloronaphthalene	U	330	1.0	330	330
2-Nitroaniline	U	820	1.0	820	820
Dimethyl Phthalate	U	330	1.0	330	330
2,6-Dinitrotoluene	U	330	1.0	330	330
Acenaphthylene	U	330	1.0	330	330
3-Nitroaniline	U	820	1.0	820	820
Acenaphthene	U	330	1.0	330	330
2,4-Dinitrophenol	U	820	1.0	820	820
Dibenzofuran	U	330	1.0	330	330
4-Nitrophenol	U	820	1.0	820	820
2,4-Dinitrotoluene	U	820	1.0	820	820
Diethylphthalate	U	330	1.0	330	330
Fluorene	U	330	1.0	330	330
4-Chlorophenyl-phenylether	U	330	1.0	330	330
4-Nitroaniline	U	820	1.0	820	820
4,6-Dinitro-2-Methylphenol	U	820	1.0	820	820
N-Nitrosodiphenylamine	U	330	1.0	330	330

KATAHDIN ANALYTICAL SERVICES
LAB CONTROL SAMPLE

Client:
Project: Old Town
PO No:
Sample Date:
Received Date:
Extraction Date: 05/31/06
Analysis Date: 06/01/06
Report Date: 06/06/2006
Matrix: SOIL

Lab ID: WG28621-2 & WG28621-3
Client ID: WG28621-LCS & WG28621-LCSD
SDG: WW2637
Extracted by: KM
Extraction Method: SW846 3550
Analyst: LRS
Analysis Method: SW846 8270C
Lab Prep Batch: WG28621
Units: ug/Kg

COMPOUND	LCS SPIKE	LCSD SPIKE	SAMPLE CONC.	LCS CONC.	LCSD CONC.	LCS %REC.		LCSD %REC.		%RPD	%RPD LIMIT	QC. LIMITS
Phenol	3333	3333	NA	2240	2320	67		70		4	50	38-125
Bis(2-Chloroethyl)ether	1667	1667	NA	986	1030	59		62		4	50	30-114
2-Chlorophenol	3333	3333	NA	2190	2280	66		68		4	50	29-130
1,3-Dichlorobenzene	1667	1667	NA	834	981	50		59		16	50	26-114
1,4-Dichlorobenzene	1667	1667	NA	853	996	51		60		15	50	21-123
1,2-Dichlorobenzene	1667	1667	NA	934	1070	56		64		14	50	28-122
2-Methylphenol	3333	3333	NA	2300	2460	69		74		7	50	6-134
2,2'-Oxybis(1-Chloropropane)	1667	1667	NA	1040	1080	62		65		4	50	32-138
N-Nitroso-di-n-propylamine	1667	1667	NA	1090	1070	65		64		2	50	35-113
3&4-Methylphenol	3333	3333	NA	2350	2560	70		77		8	50	42-114
Hexachloroethane	1667	1667	NA	884	1030	53		62		15	50	9-113
Nitrobenzene	1667	1667	NA	995	1050	60		63		5	50	37-117
Isophorone	1667	1667	NA	1040	1100	62		66		6	50	45-109
2-Nitrophenol	3333	3333	NA	2080	2270	62		68		9	50	6-130
2,4-Dimethylphenol	3333	3333	NA	2350	2690	70		81		13	50	46- 99
Bis(2-Chloroethoxy)methane	1667	1667	NA	1060	1160	64		70		9	50	38-116
2,4-Dichlorophenol	3333	3333	NA	2300	2640	69		79		14	50	42-111
1,2,4-Trichlorobenzene	1667	1667	NA	970	1070	58		64		10	50	24-131
Naphthalene	1667	1667	NA	1020	1120	61		67		9	50	36-111
4-Chloroaniline	1667	1667	NA	1190	2140	71	*	128	*	57	50	10- 95
Hexachlorobutadiene	1667	1667	NA	968	1110	58		67		14	50	20-133
4-Chloro-3-Methylphenol	3333	3333	NA	2880	3390	86		102		16	50	50-124
2-Methylnaphthalene	1667	1667	NA	1120	1220	67		73		8	50	43-123
Hexachlorocyclopentadiene	1667	1667	NA	568	678	34		41		18	50	15-136
2,4,6-Trichlorophenol	3333	3333	NA	2670	3110	80		93		15	50	45-110
2,4,5-Trichlorophenol	3333	3333	NA	2240	2610	67		78		15	50	50-110
2-Chloronaphthalene	1667	1667	NA	1040	1160	62		70		11	50	39- 99
2-Nitroaniline	1667	1667	NA	1260	1390	76		83		10	50	43-131
Dimethyl Phthalate	1667	1667	NA	1290	1460	77		88		12	50	61-112
2,6-Dinitrotoluene	1667	1667	NA	1190	1310	71		79		10	50	58-106
Acenaphthylene	1667	1667	NA	1160	1320	70		79		13	50	50-115
3-Nitroaniline	1667	1667	NA	139	111	8	*	7	*	22	50	25-110
Acenaphthene	1667	1667	NA	1120	1240	67		74		10	50	50-118
2,4-Dinitrophenol	3333	3333	NA	1060	748	32		22		34	50	3- 93
Dibenzofuran	1667	1667	NA	1210	1340	73		80		10	50	53-128
4-Nitrophenol	3333	3333	NA	2580	2840	77		85		10	50	28-137
2,4-Dinitrotoluene	1667	1667	NA	1290	1400	77		84		8	50	42-120
Diethylphthalate	1667	1667	NA	1270	1400	76		84		10	50	57-108
Fluorene	1667	1667	NA	1200	1340	72		80		11	50	63-105
4-Chlorophenyl-phenylether	1667	1667	NA	1180	1300	71		78		10	50	63-110
4-Nitroaniline	1667	1667	NA	1310	1390	79		83		6	50	22- 93
4,6-Dinitro-2-Methylphenol	3333	3333	NA	1840	1620	55		49		13	50	27-111
N-Nitrosodiphenylamine	1667	1667	NA	1230	1390	74		83		12	50	50-115
4-Bromophenyl-phenylether	1667	1667	NA	1260	1440	76		86		13	50	48-133
Hexachlorobenzene	1667	1667	NA	1240	1420	74		85		14	50	33-154

KATAHDIN ANALYTICAL SERVICES
LAB CONTROL SAMPLE

Client:
Project: Old Town
PO No:
Sample Date:
Received Date:
Extraction Date: 05/31/06
Analysis Date: 06/01/06
Report Date: 06/06/2006
Matrix: SOIL

Lab ID: WG28621-2 & WG28621-3
Client ID: WG28621-LCS & WG28621-LCSD
SDG: WW2637
Extracted by: KM
Extraction Method: SW846 3550
Analyst: LRS
Analysis Method: SW846 8270C
Lab Prep Batch: WG28621
Units: ug/Kg

COMPOUND	LCS SPIKE	LCSD SPIKE	SAMPLE CONC.	LCS CONC.	LCSD CONC.	LCS %REC.	LCSD %REC.	%RPD	%RPD LIMIT	QC. LIMITS
Pentachlorophenol	3333	3333	NA	2700	3040	81	91	12	50	32-139
Phenanthrene	1667	1667	NA	1240	1380	74	83	11	50	55-121
Anthracene	1667	1667	NA	1220	1410	73	85	14	50	60-122
Carbazole	1667	1667	NA	1410	1500	85	90	6	50	51-119
Di-n-butylphthalate	1667	1667	NA	1370	1470	82	88	7	50	40-120
Fluoranthene	1667	1667	NA	1380	1480	83	89	7	50	41-135
Pyrene	1667	1667	NA	1340	1570	80	94	16	50	50-114
Butylbenzylphthalate	1667	1667	NA	1410	1600	85	96	13	50	46-115
Benzo(a)anthracene	1667	1667	NA	1310	1420	79	85	8	50	45-111
3,3'-Dichlorobenzidine	1667	1667	NA	3520	4660	* 211	* 280	28	50	10-130
Chrysene	1667	1667	NA	1360	1520	82	91	11	50	45-124
bis(2-Ethylhexyl)phthalate	1667	1667	NA	1420	1620	85	97	13	50	45-125
Di-n-octylphthalate	1667	1667	NA	1510	1860	91	112	21	50	39-126
Benzo(b)fluoranthene	1667	1667	NA	1260	1400	76	84	10	50	47-114
Benzo(k)fluoranthene	1667	1667	NA	1380	1570	83	94	13	50	46-125
Benzo(a)pyrene	1667	1667	NA	1300	1420	78	85	9	50	45-119
Indeno(1,2,3-cd)pyrene	1667	1667	NA	1220	1180	73	71	3	50	38-131
Dibenzo(a,h)anthracene	1667	1667	NA	1230	1230	74	74	0.0	50	37-125
Benzo(g,h,i)perylene	1667	1667	NA	1180	1200	71	72	2	50	37-133

FORM III SV-2

R1172.D & R1173.D

FORM 4

SEMIVOLATILE METHOD BLANK SUMMARY

CLIENT SAMPLE ID

| WG28661-BLANK |

Lab Name: KATAHDIN ANALYTICAL SERVICES Lab Code: KAS

Project: OLD TOWN SDG No.: WW2637

Lab File ID: R1174 Lab Sample ID: WG28661-1

Instrument ID: GCMS-R Date Extracted: 05/31/06

Matrix: (soil/water) WATER Date Analyzed: 06/01/06

Level: (low/med) LOW Time Analyzed: 1344

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES, MS and MSD:

	CLIENT SAMPLE ID	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED	TIME ANALYZED
01	WG28661-LCS	WG28661-2	R1175	06/01/06	1430
02	SB-25	WW2637-1	R1176	06/01/06	1517
03	MW-1	WW2637-2	R1177	06/01/06	1604
04	MW-2	WW2637-3	R1178	06/01/06	1651
05	SB-26	WW2637-4	R1180	06/02/06	0815
06	SB-100	WW2637-5	R1181	06/02/06	0859
07	SB-25MS	WG28661-3	R1184	06/02/06	1116
08	SB-25MSD	WG28661-4	R1185	06/02/06	1201
09					
10					
11					
12					
13					
14					
15					
16					
17					
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21					
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25					
26					
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28					
29					
30					

COMMENTS:

KATAHDIN ANALYTICAL SERVICES
Report of Analytical Results

Client:
Project: Old Town
PO No:
Sample Date:
Received Date:
Extraction Date: 05/31/06
Analysis Date: 01-JUN-2006 13:44
Report Date: 06/06/2006
Matrix: WATER
% Solids: NA

Lab ID: WG28661-1
Client ID: WG28661-Blank
SDG: WW2637
Extracted by: GN
Extraction Method: SW846 3510
Analyst: LRS
Analysis Method: SW846 8270C
Lab Prep Batch: WG28661
Units: ug/L

Compound	Flags	Results	DF	PQL	Adj.PQL
Phenol	U	10	1.0	10	10
Bis(2-Chloroethyl)ether	U	10	1.0	10	10
2-Chlorophenol	U	10	1.0	10	10
1,3-Dichlorobenzene	U	10	1.0	10	10
1,4-Dichlorobenzene	U	10	1.0	10	10
1,2-Dichlorobenzene	U	10	1.0	10	10
2-Methylphenol	U	10	1.0	10	10
2,2'-Oxybis(1-Chloropropane)	U	10	1.0	10	10
N-Nitroso-di-n-propylamine	U	10	1.0	10	10
3&4-Methylphenol	U	10	1.0	10	10
Hexachloroethane	U	10	1.0	10	10
Nitrobenzene	U	10	1.0	10	10
Isophorone	U	10	1.0	10	10
2-Nitrophenol	U	10	1.0	10	10
2,4-Dimethylphenol	U	10	1.0	10	10
Bis(2-Chloroethoxy)methane	U	10	1.0	10	10
2,4-Dichlorophenol	U	10	1.0	10	10
1,2,4-Trichlorobenzene	U	10	1.0	10	10
Naphthalene	U	10	1.0	10	10
4-Chloroaniline	U	10	1.0	10	10
Hexachlorobutadiene	U	10	1.0	10	10
4-Chloro-3-Methylphenol	U	10	1.0	10	10
2-Methylnaphthalene	U	10	1.0	10	10
Hexachlorocyclopentadiene	U	10	1.0	10	10
2,4,6-Trichlorophenol	U	10	1.0	10	10
2,4,5-Trichlorophenol	U	25	1.0	25	25
2-Chloronaphthalene	U	10	1.0	10	10
2-Nitroaniline	U	25	1.0	25	25
Dimethyl Phthalate	U	10	1.0	10	10
2,6-Dinitrotoluene	U	10	1.0	10	10
Acenaphthylene	U	10	1.0	10	10
3-Nitroaniline	U	25	1.0	25	25
Acenaphthene	U	10	1.0	10	10
2,4-Dinitrophenol	U	25	1.0	25	25
Dibenzofuran	U	10	1.0	10	10
4-Nitrophenol	U	25	1.0	25	25
2,4-Dinitrotoluene	U	10	1.0	10	10
Diethylphthalate	U	10	1.0	10	10
Fluorene	U	10	1.0	10	10
4-Chlorophenyl-phenylether	U	10	1.0	10	10
4-Nitroaniline	U	25	1.0	25	25
4,6-Dinitro-2-Methylphenol	U	25	1.0	25	25
N-Nitrosodiphenylamine	U	10	1.0	10	10

KATAHDIN ANALYTICAL SERVICES
Report of Analytical Results

Client:
Project: Old Town
PO No:
Sample Date:
Received Date:
Extraction Date: 05/31/06
Analysis Date: 01-JUN-2006 13:44
Report Date: 06/06/2006
Matrix: WATER
% Solids: NA

Lab ID: WG28661-1
Client ID: WG28661-Blank
SDG: WW2637
Extracted by: GN
Extraction Method: SW846 3510
Analyst: LRS
Analysis Method: SW846 8270C
Lab Prep Batch: WG28661
Units: ug/L

Compound	Flags	Results	DF	PQL	Adj.PQL
4-Bromophenyl-phenylether	U	10	1.0	10	10
Hexachlorobenzene	U	10	1.0	10	10
Pentachlorophenol	U	25	1.0	25	25
Phenanthrene	U	10	1.0	10	10
Anthracene	U	10	1.0	10	10
Carbazole	U	10	1.0	10	10
Di-n-butylphthalate	U	10	1.0	10	10
Fluoranthene	U	10	1.0	10	10
Pyrene	U	10	1.0	10	10
Butylbenzylphthalate	U	10	1.0	10	10
Benzo(a)anthracene	U	10	1.0	10	10
3,3'-Dichlorobenzidine	U	10	1.0	10	10
Chrysene	U	10	1.0	10	10
bis(2-Ethylhexyl)phthalate	U	10	1.0	10	10
Di-n-octylphthalate	U	10	1.0	10	10
Benzo(b)fluoranthene	U	10	1.0	10	10
Benzo(k)fluoranthene	U	10	1.0	10	10
Benzo(a)pyrene	U	10	1.0	10	10
Indeno(1,2,3-cd)pyrene	U	10	1.0	10	10
Dibenzo(a,h)anthracene	U	10	1.0	10	10
Benzo(g,h,i)perylene	U	10	1.0	10	10
2-Fluorophenol		41%			
Phenol-D6		32%			
Nitrobenzene-D5		* 50%			
2-Fluorobiphenyl		* 56%			
2,4,6-Tribromophenol		78%			
Terphenyl-D14		90%			

KATAHDIN ANALYTICAL SERVICES
LAB CONTROL SAMPLE

Client:
Project: Old Town
PO No:
Sample Date:
Received Date:
Extraction Date: 05/31/06
Analysis Date: 06/01/06
Report Date: 06/06/2006
Matrix: WATER

Lab ID: WG28661-2
Client ID: WG28661-LCS
SDG: WW2637
Extracted by: GN
Extraction Method: SW846 3510
Analyst: LRS
Analysis Method: SW846 8270C
Lab Prep Batch: WG28661
Units: ug/L

COMPOUND	LCS SPIKE	SAMPLE CONC.	LCS CONC.		%REC.	QC. LIMITS
Phenol	100	NA	42		42	0-126
Bis(2-Chloroethyl)ether	50	NA	38		76	35-109
2-Chlorophenol	100	NA	82		82	52-112
1,3-Dichlorobenzene	50	NA	32		64	20-100
1,4-Dichlorobenzene	50	NA	32		64	19-103
1,2-Dichlorobenzene	50	NA	35		70	25- 93
2-Methylphenol	100	NA	78		78	31-105
2,2'-Oxybis(1-Chloropropane)	50	NA	40		80	32-108
N-Nitroso-di-n-propylamine	50	NA	40		80	36-106
3&4-Methylphenol	100	NA	70		70	24-116
Hexachloroethane	50	NA	34		68	30- 95
Nitrobenzene	50	NA	34		69	44-112
Isophorone	50	NA	37		74	50-110
2-Nitrophenol	100	NA	82		82	49-108
2,4-Dimethylphenol	100	NA	88		88	48-105
Bis(2-Chloroethoxy)methane	50	NA	38		75	52-101
2,4-Dichlorophenol	100	NA	87		87	57-110
1,2,4-Trichlorobenzene	50	NA	33		67	31-114
Naphthalene	50	NA	36		71	33-109
4-Chloroaniline	50	NA	95	*	190	29- 98
Hexachlorobutadiene	50	NA	34		67	30-107
4-Chloro-3-Methylphenol	100	NA	101		101	46-150
2-Methylnaphthalene	50	NA	39		78	48-121
Hexachlorocyclopentadiene	50	NA	19		39	0-106
2,4,6-Trichlorophenol	100	NA	95		95	69-105
2,4,5-Trichlorophenol	100	NA	81		81	58-117
2-Chloronaphthalene	50	NA	35		70	33-119
2-Nitroaniline	50	NA	41		81	45-127
Dimethyl Phthalate	50	NA	29		57	4-149
2,6-Dinitrotoluene	50	NA	38		76	51-120
Acenaphthylene	50	NA	41		82	49-117
3-Nitroaniline	50	NA	4.1	*	8	20-125
Acenaphthene	50	NA	38		76	45-118
2,4-Dinitrophenol	100	NA	59		59	15-140
Dibenzofuran	50	NA	39		79	53-124
4-Nitrophenol	100	NA	40		40	0-158
2,4-Dinitrotoluene	50	NA	40		80	44-133
Diethylphthalate	50	NA	37		73	20-139
Fluorene	50	NA	39		78	46-127
4-Chlorophenyl-phenylether	50	NA	38		77	48-136
4-Nitroaniline	50	NA	35		69	15-145
4,6-Dinitro-2-Methylphenol	100	NA	83		83	4-150
N-Nitrosodiphenylamine	50	NA	40		80	30-150
4-Bromophenyl-phenylether	50	NA	41		82	68-129
Hexachlorobenzene	50	NA	42		84	67-141

KATAHDIN ANALYTICAL SERVICES
LAB CONTROL SAMPLE

Client:
Project: Old Town
PO No:
Sample Date:
Received Date:
Extraction Date: 05/31/06
Analysis Date: 06/01/06
Report Date: 06/06/2006
Matrix: WATER

Lab ID: WG28661-2
Client ID: WG28661-LCS
SDG: WW2637
Extracted by: GN
Extraction Method: SW846 3510
Analyst: LRS
Analysis Method: SW846 8270C
Lab Prep Batch: WG28661
Units: ug/L

COMPOUND	LCS SPIKE	SAMPLE CONC.	LCS CONC.		%REC.	QC. LIMITS
Pentachlorophenol	100	NA	106		106	10-174
Phenanthrene	50	NA	41		82	51-139
Anthracene	50	NA	42		85	53-133
Carbazole	50	NA	40		81	35-157
Di-n-butylphthalate	50	NA	43		86	14-165
Fluoranthene	50	NA	45		90	32-172
Pyrene	50	NA	47		93	49-119
Butylbenzylphthalate	50	NA	46		91	30-126
Benzo(a)anthracene	50	NA	42		83	53-111
3,3'-Dichlorobenzidine	50	NA	125	*	250	38-101
Chrysene	50	NA	47		95	59-120
bis(2-Ethylhexyl)phthalate	50	NA	49		98	39-125
Di-n-octylphthalate	50	NA	53		106	31-136
Benzo(b)fluoranthene	50	NA	40		80	52-120
Benzo(k)fluoranthene	50	NA	47		95	56-126
Benzo(a)pyrene	50	NA	42		84	51-120
Indeno(1,2,3-cd)pyrene	50	NA	36		71	32-143
Dibenzo(a,h)anthracene	50	NA	38		77	37-134
Benzo(g,h,i)perylene	50	NA	38		77	32-143

FORM III SV-1

R1175.D

KATAHDIN ANALYTICAL SERVICES
MATRIX SPIKE/MATRIX SPIKE DUPLICATE RECOVERY

Client: Woodard & Curran
Project: Old Town
PO No:
Sample Date: 05/30/06
Received Date: 05/31/06
Extraction Date: 05/31/06
Analysis Date: 06/02/06
Report Date: 06/07/2006
Matrix: WATER

Lab ID: WG28661-3 & WG28661-4
Client ID: SB-25MS & SB-25MSD
SDG: WW2637
Extracted by: GN
Extraction Method: SW846 3510
Analyst: LRS
Analysis Method: SW846 8270C
Lab Prep Batch: WG28661
Units: ug/L

COMPOUND	MS SPIKE	MSD SPIKE	SAMPLE CONC.	MS CONC.	MSD CONC.	MS %REC.	MSD %REC.	%RPD	%RPD LIMIT	QC. LIMITS
Phenol	100	100	0.00	35	44	35	44	21	30	0-126
Bis(2-Chloroethyl)ether	50	50	0.00	34	42	68	85	22	30	35-109
2-Chlorophenol	100	100	0.00	67	83	67	83	21	30	52-112
1,3-Dichlorobenzene	50	50	0.00	29	35	57	70	20	30	20-100
1,4-Dichlorobenzene	50	50	0.00	28	35	57	70	20	30	19-103
1,2-Dichlorobenzene	50	50	0.00	32	39	64	77	19	30	25- 93
2-Methylphenol	100	100	0.00	62	77	62	77	20	30	31-105
2,2'-Oxybis(1-Chloropropane)	50	50	0.00	38	45	75	89	17	30	32-108
N-Nitroso-di-n-propylamine	50	50	0.00	38	44	75	88	16	30	36-106
3&4-Methylphenol	100	100	0.00	52	65	52	65	22	30	24-116
Hexachloroethane	50	50	0.00	28	34	56	68	20	30	30- 95
Nitrobenzene	50	50	0.00	32	38	63	77	19	30	44-112
Isophorone	50	50	0.00	34	41	69	82	17	30	50-110
2-Nitrophenol	100	100	0.00	71	88	71	88	21	30	49-108
2,4-Dimethylphenol	100	100	0.00	59	76	59	76	26	30	48-105
Bis(2-Chloroethoxy)methane	50	50	0.00	35	41	71	83	15	30	52-101
2,4-Dichlorophenol	100	100	0.00	67	83	67	83	21	30	57-110
1,2,4-Trichlorobenzene	50	50	0.00	30	36	59	72	20	30	31-114
Naphthalene	50	50	0.00	32	40	65	79	20	30	33-109
4-Chloroaniline	50	50	0.00	32	38	64	77	19	30	29- 98
Hexachlorobutadiene	50	50	0.00	30	36	61	71	16	30	30-107
4-Chloro-3-Methylphenol	100	100	0.00	81	96	81	96	17	30	46-150
2-Methylnaphthalene	50	50	0.00	35	42	70	84	18	30	48-121
Hexachlorocyclopentadiene	50	50	0.00	7.8	9.2	16	18	17	30	0-106
2,4,6-Trichlorophenol	100	100	0.00	78	91	78	91	15	30	69-105
2,4,5-Trichlorophenol	100	100	0.00	69	79	69	79	14	30	58-117
2-Chloronaphthalene	50	50	0.00	32	38	63	76	18	30	33-119
2-Nitroaniline	50	50	0.00	39	44	79	88	11	30	45-127
Dimethyl Phthalate	50	50	0.00	38	42	76	84	10	30	4-149
2,6-Dinitrotoluene	50	50	0.00	36	40	73	80	9	30	51-120
Acenaphthylene	50	50	0.00	37	43	75	86	14	30	49-117
3-Nitroaniline	50	50	0.00	0.0	4.1	* 0	* 8		30	20-125
Acenaphthene	50	50	0.00	35	40	70	81	15	30	45-118
2,4-Dinitrophenol	100	100	0.00	44	53	44	53	17	30	15-140
Dibenzofuran	50	50	0.00	36	42	73	83	13	30	53-124
4-Nitrophenol	100	100	0.00	43	44	43	44	3	30	0-158
2,4-Dinitrotoluene	50	50	0.00	38	41	76	82	6	30	44-133
Diethylphthalate	50	50	0.00	39	41	78	83	6	30	20-139
Fluorene	50	50	0.00	37	40	74	79	7	30	46-127
4-Chlorophenyl-phenylether	50	50	0.00	35	39	71	78	10	30	48-136
4-Nitroaniline	50	50	0.00	39	41	78	82	5	30	15-145
4,6-Dinitro-2-Methylphenol	100	100	0.00	59	68	59	68	14	30	4-150
N-Nitrosodiphenylamine	50	50	0.00	40	43	81	85	5	30	30-150
4-Bromophenyl-phenylether	50	50	0.00	40	43	81	86	6	30	68-129
Hexachlorobenzene	50	50	0.00	40	42	80	84	5	30	67-141

KATAHDIN ANALYTICAL SERVICES
MATRIX SPIKE/MATRIX SPIKE DUPLICATE RECOVERY

Client: Woodard & Curran
Project: Old Town
PO No:
Sample Date: 05/30/06
Received Date: 05/31/06
Extraction Date: 05/31/06
Analysis Date: 06/02/06
Report Date: 06/07/2006
Matrix: WATER

Lab ID: WG28661-3 & WG28661-4
Client ID: SB-25MS & SB-25MSD
SDG: WW2637
Extracted by: GN
Extraction Method: SW846 3510
Analyst: LRS
Analysis Method: SW846 8270C
Lab Prep Batch: WG28661
Units: ug/L

COMPOUND	MS SPIKE	MSD SPIKE	SAMPLE CONC.	MS CONC.	MSD CONC.	MS %REC.	MSD %REC.	%RPD	%RPD LIMIT	QC. LIMITS
Pentachlorophenol	100	100	0.00	91	98	91	98	7	30	10-174
Phenanthrene	50	50	0.00	39	41	79	83	5	30	51-139
Anthracene	50	50	0.00	40	42	81	84	4	30	53-133
Carbazole	50	50	0.00	42	44	85	89	5	30	35-157
Di-n-butylphthalate	50	50	0.00	42	44	84	87	4	30	14-165
Fluoranthene	50	50	0.00	37	39	74	78	6	30	32-172
Pyrene	50	50	0.00	56	59	113	118	5	30	49-119
Butylbenzylphthalate	50	50	0.00	58	61	117	121	4	30	30-126
Benzo(a)anthracene	50	50	0.00	39	40	77	80	4	30	53-111
3,3'-Dichlorobenzidine	50	50	0.00	105	100	* 210	* 200	5	30	38-101
Chrysene	50	50	0.00	43	46	86	92	6	30	59-120
bis(2-Ethylhexyl)phthalate	50	50	0.00	55	58	109	116	6	30	39-125
Di-n-octylphthalate	50	50	0.00	66	72	132	* 143	8	30	31-136
Benzo(b)fluoranthene	50	50	0.00	37	40	75	80	7	30	52-120
Benzo(k)fluoranthene	50	50	0.00	46	50	92	100	8	30	56-126
Benzo(a)pyrene	50	50	0.00	36	44	72	89	21	30	51-120
Indeno(1,2,3-cd)pyrene	50	50	0.00	32	36	64	73	12	30	32-143
Dibenzo(a,h)anthracene	50	50	0.00	36	42	72	84	14	30	37-134
Benzo(g,h,i)perylene	50	50	0.00	32	36	65	72	10	30	32-143

FORM III SV-1

R1184.D & R1185.D

FORM 4
SEMIVOLATILE METHOD BLANK SUMMARY

CLIENT SAMPLE ID

WG28761-BLANK

Lab Name: KATAHDIN ANALYTICAL SERVICES Lab Code: KAS

Project: OLD TOWN SDG No.: WW2637

Lab File ID: X0016 Lab Sample ID: WG28761-1

Instrument ID: GCMS-X Date Extracted: 06/04/06

Matrix: (soil/water) SOIL Date Analyzed: 06/06/06

Level: (low/med) LOW Time Analyzed: 1123

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES, MS and MSD:

	CLIENT SAMPLE ID	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED	TIME ANALYZED
01	WG28761-LCS	WG28761-2	X0017	06/06/06	1208
02	SS05MS	WG28761-3	X0018	06/06/06	1253
03	SS05MSD	WG28761-4	X0019	06/06/06	1338
04					
05					
06					
07					
08					
09					
10					
11					
12					
13					
14					
15					
16					
17					
18					
19					
20					
21					
22					
23					
24					
25					
26					
27					
28					
29					
30					

COMMENTS:

KATAHDIN ANALYTICAL SERVICES
Report of Analytical Results

Client:
Project: Old Town
PO No:
Sample Date:
Received Date:
Extraction Date: 06/04/06
Analysis Date: 06-JUN-2006 11:23
Report Date: 06/06/2006
Matrix: SOIL
% Solids: 100

Lab ID: WG28761-1
Client ID: WG28761-Blank
SDG: WW2637
Extracted by: GN
Extraction Method: SW846 3550
Analyst: JCG
Analysis Method: SW846 8270C
Lab Prep Batch: WG28761
Units: ug/Kg

Compound	Flags	Results	DF	PQL	Adj.PQL
Phenol	U	330	1.0	330	330
Bis(2-Chloroethyl)ether	U	330	1.0	330	330
2-Chlorophenol	U	330	1.0	330	330
1,3-Dichlorobenzene	U	330	1.0	330	330
1,4-Dichlorobenzene	U	330	1.0	330	330
1,2-Dichlorobenzene	U	330	1.0	330	330
2-Methylphenol	U	330	1.0	330	330
2,2'-Oxybis(1-Chloropropane)	U	330	1.0	330	330
N-Nitroso-di-n-propylamine	U	330	1.0	330	330
3&4-Methylphenol	U	330	1.0	330	330
Hexachloroethane	U	330	1.0	330	330
Nitrobenzene	U	330	1.0	330	330
Isophorone	U	330	1.0	330	330
2-Nitrophenol	U	330	1.0	330	330
2,4-Dimethylphenol	U	330	1.0	330	330
Bis(2-Chloroethoxy)methane	U	330	1.0	330	330
2,4-Dichlorophenol	U	330	1.0	330	330
1,2,4-Trichlorobenzene	U	330	1.0	330	330
Naphthalene	U	330	1.0	330	330
4-Chloroaniline	U	330	1.0	330	330
Hexachlorobutadiene	U	330	1.0	330	330
4-Chloro-3-Methylphenol	U	330	1.0	330	330
2-Methylnaphthalene	U	330	1.0	330	330
Hexachlorocyclopentadiene	U	330	1.0	330	330
2,4,6-Trichlorophenol	U	330	1.0	330	330
2,4,5-Trichlorophenol	U	820	1.0	820	820
2-Chloronaphthalene	U	330	1.0	330	330
2-Nitroaniline	U	820	1.0	820	820
Dimethyl Phthalate	U	330	1.0	330	330
2,6-Dinitrotoluene	U	330	1.0	330	330
Acenaphthylene	U	330	1.0	330	330
3-Nitroaniline	U	820	1.0	820	820
Acenaphthene	U	330	1.0	330	330
2,4-Dinitrophenol	U	820	1.0	820	820
Dibenzofuran	U	330	1.0	330	330
4-Nitrophenol	U	820	1.0	820	820
2,4-Dinitrotoluene	U	820	1.0	820	820
Diethylphthalate	U	330	1.0	330	330
Fluorene	U	330	1.0	330	330
4-Chlorophenyl-phenylether	U	330	1.0	330	330
4-Nitroaniline	U	820	1.0	820	820
4,6-Dinitro-2-Methylphenol	U	820	1.0	820	820
N-Nitrosodiphenylamine	U	330	1.0	330	330

KATAHDIN ANALYTICAL SERVICES
Report of Analytical Results

Client:
Project: Old Town
PO No:
Sample Date:
Received Date:
Extraction Date: 06/04/06
Analysis Date: 06-JUN-2006 11:23
Report Date: 06/06/2006
Matrix: SOIL
% Solids: 100

Lab ID: WG28761-1
Client ID: WG28761-Blank
SDG: WW2637
Extracted by: GN
Extraction Method: SW846 3550
Analyst: JCG
Analysis Method: SW846 8270C
Lab Prep Batch: WG28761
Units: ug/Kg

Compound	Flags	Results	DF	PQL	Adj.PQL
4-Bromophenyl-phenylether	U	330	1.0	330	330
Hexachlorobenzene	U	330	1.0	330	330
Pentachlorophenol	U	820	1.0	820	820
Phenanthrene	U	330	1.0	330	330
Anthracene	U	330	1.0	330	330
Carbazole	U	330	1.0	330	330
Di-n-butylphthalate	U	330	1.0	330	330
Fluoranthene	U	330	1.0	330	330
Pyrene	U	330	1.0	330	330
Butylbenzylphthalate	U	330	1.0	330	330
Benzo(a)anthracene	U	330	1.0	330	330
3,3'-Dichlorobenzidine	U	330	1.0	330	330
Chrysene	U	330	1.0	330	330
bis(2-Ethylhexyl)phthalate	U	330	1.0	330	330
Di-n-octylphthalate	U	330	1.0	330	330
Benzo(b)fluoranthene	U	330	1.0	330	330
Benzo(k)fluoranthene	U	330	1.0	330	330
Benzo(a)pyrene	U	330	1.0	330	330
Indeno(1,2,3-cd)pyrene	U	330	1.0	330	330
Dibenzo(a,h)anthracene	U	330	1.0	330	330
Benzo(g,h,i)perylene	U	330	1.0	330	330
2-Fluorophenol		46%			
Phenol-D6		82%			
Nitrobenzene-D5		78%			
2-Fluorobiphenyl		90%			
2,4,6-Tribromophenol		*133%			
Terphenyl-D14		97%			

KATAHDIN ANALYTICAL SERVICES
LAB CONTROL SAMPLE

Client:
Project: Old Town
PO No:
Sample Date:
Received Date:
Extraction Date: 06/04/06
Analysis Date: 06/06/06
Report Date: 06/06/2006
Matrix: SOIL

Lab ID: WG28761-2
Client ID: WG28761-LCS
SDG: WW2637
Extracted by: GN
Extraction Method: SW846 3550
Analyst: JCG
Analysis Method: SW846 8270C
Lab Prep Batch: WG28761
Units: ug/Kg

COMPOUND	LCS SPIKE	SAMPLE CONC.	LCS CONC.		%REC.	QC. LIMITS
Phenol	3333	NA	2860		86	38-125
Bis(2-Chloroethyl)ether	1667	NA	2040	*	122	30-114
2-Chlorophenol	3333	NA	3180		95	29-130
1,3-Dichlorobenzene	1667	NA	1320		79	26-114
1,4-Dichlorobenzene	1667	NA	1410		85	21-123
1,2-Dichlorobenzene	1667	NA	1480		89	28-122
2-Methylphenol	3333	NA	3460		104	6-134
2,2'-Oxybis(1-Chloropropane)	1667	NA	1220		73	23-138
N-Nitroso-di-n-propylamine	1667	NA	1490		89	35-113
3&4-Methylphenol	3333	NA	3340		100	42-114
Hexachloroethane	1667	NA	1200		72	9-113
Nitrobenzene	1667	NA	1540		92	37-117
Isophorone	1667	NA	1530		92	45-109
2-Nitrophenol	3333	NA	3050		92	6-130
2,4-Dimethylphenol	3333	NA	3050		92	46- 99
Bis(2-Chloroethoxy)methane	1667	NA	1430		86	38-116
2,4-Dichlorophenol	3333	NA	3560		107	42-111
1,2,4-Trichlorobenzene	1667	NA	1700		102	24-131
Naphthalene	1667	NA	1550		93	36-111
4-Chloroaniline	1667	NA	792		48	10- 95
Hexachlorobutadiene	1667	NA	2060		124	20-133
4-Chloro-3-Methylphenol	3333	NA	3420		103	50-124
2-Methylnaphthalene	1667	NA	1560		94	43-123
Hexachlorocyclopentadiene	1667	NA	987		59	15-136
2,4,6-Trichlorophenol	3333	NA	3710	*	111	45-110
2,4,5-Trichlorophenol	3333	NA	3140		94	50-110
2-Chloronaphthalene	1667	NA	1280		77	39- 99
2-Nitroaniline	1667	NA	1450		87	43-131
Dimethyl Phthalate	1667	NA	1740		104	61-112
2,6-Dinitrotoluene	1667	NA	1620		97	58-106
Acenaphthylene	1667	NA	1440		86	50-115
3-Nitroaniline	1667	NA	1100		66	25-110
Acenaphthene	1667	NA	1380		83	50-118
2,4-Dinitrophenol	3333	NA	1130		34	3- 93
Dibenzofuran	1667	NA	1580		95	53-128
4-Nitrophenol	3333	NA	3880		116	28-137
2,4-Dinitrotoluene	1667	NA	1540		92	42-120
Diethylphthalate	1667	NA	1590		95	57-108
Fluorene	1667	NA	1530		92	63-105
4-Chlorophenyl-phenylether	1667	NA	1760		106	63-110
4-Nitroaniline	1667	NA	950		57	22- 93
4,6-Dinitro-2-Methylphenol	3333	NA	2350		70	27-111
N-Nitrosodiphenylamine	1667	NA	1480		89	50-115
4-Bromophenyl-phenylether	1667	NA	2310	*	139	48-133
Hexachlorobenzene	1667	NA	2810	*	169	33-154

KATAHDIN ANALYTICAL SERVICES
LAB CONTROL SAMPLE

Client:
Project: Old Town
PO No:
Sample Date:
Received Date:
Extraction Date: 06/04/06
Analysis Date: 06/06/06
Report Date: 06/06/2006
Matrix: SOIL

Lab ID: WG28761-2
Client ID: WG28761-LCS
SDG: WW2637
Extracted by: GN
Extraction Method: SW846 3550
Analyst: JCG
Analysis Method: SW846 8270C
Lab Prep Batch: WG28761
Units: ug/Kg

COMPOUND	LCS SPIKE	SAMPLE CONC.	LCS CONC.		%REC.	QC. LIMITS
Pentachlorophenol	3333	NA	4420		133	32-139
Phenanthrene	1667	NA	1880		113	55-121
Anthracene	1667	NA	1910		115	60-122
Carbazole	1667	NA	1600		96	51-119
Di-n-butylphthalate	1667	NA	1720		103	40-120
Fluoranthene	1667	NA	1940		116	41-135
Pyrene	1667	NA	2170	*	130	50-114
Butylbenzylphthalate	1667	NA	1520		91	46-115
Benzo(a)anthracene	1667	NA	1840		110	45-111
3,3'-Dichlorobenzidine	1667	NA	1350		81	10-130
Chrysene	1667	NA	1880		113	45-124
bis(2-Ethylhexyl)phthalate	1667	NA	1610		97	45-125
Di-n-octylphthalate	1667	NA	1340		80	39-126
Benzo(b)fluoranthene	1667	NA	1610		97	47-114
Benzo(k)fluoranthene	1667	NA	1660		100	46-125
Benzo(a)pyrene	1667	NA	1640		98	45-119
Indeno(1,2,3-cd)pyrene	1667	NA	1780		107	38-131
Dibenzo(a,h)anthracene	1667	NA	1710		103	37-125
Benzo(g,h,i)perylene	1667	NA	1700		102	37-133

FORM III SV-2

X0017.D

KATAHDIN ANALYTICAL SERVICES
MATRIX SPIKE/MATRIX SPIKE DUPLICATE RECOVERY

Client: Woodard & Curran
Project: Old Town
PO No:
Sample Date: 05/30/06
Received Date: 05/31/06
Extraction Date: 06/04/06
Analysis Date: 06/06/06
Report Date: 06/06/2006
Matrix: SOIL

Lab ID: WG28761-3 & WG28761-4
Client ID: SS05MS & SS05MSD
SDG: WW2637
Extracted by: GN
Extraction Method: SW846 3550
Analyst: JCG
Analysis Method: SW846 8270C
Lab Prep Batch: WG28761
Units: ug/Kg

COMPOUND	MS SPIKE	MSD SPIKE	SAMPLE CONC.	MS CONC.	MSD CONC.	MS %REC.	MSD %REC.	%RPD	%RPD LIMIT	QC. LIMITS
Phenol	5556	5556	0.00	4440	3840	80	69	14	50	38-125
Bis(2-Chloroethyl)ether	2778	2778	0.00	2230	2020	80	73	10	50	30-114
2-Chlorophenol	5556	5556	0.00	4610	3840	83	69	18	50	29-130
1,3-Dichlorobenzene	2778	2778	0.00	1700	1530	61	55	10	50	26-114
1,4-Dichlorobenzene	2778	2778	0.00	1880	1600	68	58	16	50	21-123
1,2-Dichlorobenzene	2778	2778	0.00	1980	1760	71	63	12	50	28-122
2-Methylphenol	5556	5556	0.00	4500	3970	81	71	12	50	6-134
2,2'-Oxybis(1-Chloropropane)	2778	2778	0.00	1810	1520	65	55	17	50	23-138
N-Nitroso-di-n-propylamine	2778	2778	0.00	2190	2070	79	74	6	50	35-113
3&4-Methylphenol	5556	5556	0.00	4690	4220	84	76	10	50	42-114
Hexachloroethane	2778	2778	0.00	1590	1420	57	51	11	50	9-113
Nitrobenzene	2778	2778	0.00	2300	1920	83	69	18	50	37-117
Isophorone	2778	2778	0.00	2380	1940	86	70	20	50	45-109
2-Nitrophenol	5556	5556	0.00	5270	4100	95	74	25	50	6-130
2,4-Dimethylphenol	5556	5556	0.00	4000	3760	72	68	6	50	46- 99
Bis(2-Chloroethoxy)methane	2778	2778	0.00	2260	1940	81	70	15	50	38-116
2,4-Dichlorophenol	5556	5556	0.00	5390	4780	97	86	12	50	42-111
1,2,4-Trichlorobenzene	2778	2778	0.00	2450	2100	88	76	15	50	24-131
Naphthalene	2778	2778	0.00	2300	1870	83	67	21	50	36-111
4-Chloroaniline	2778	2778	0.00	992	902	36	32	10	50	10- 95
Hexachlorobutadiene	2778	2778	0.00	3110	2750	112	99	12	50	20-133
4-Chloro-3-Methylphenol	5556	5556	0.00	5110	4560	92	82	11	50	50-124
2-Methylnaphthalene	2778	2778	0.00	2420	2070	87	74	16	50	43-123
Hexachlorocyclopentadiene	2778	2778	0.00	839	751	30	27	11	50	15-136
2,4,6-Trichlorophenol	5556	5556	0.00	5890	4930	106	89	18	50	45-110
2,4,5-Trichlorophenol	5556	5556	0.00	5200	4700	94	84	10	50	50-110
2-Chloronaphthalene	2778	2778	0.00	2080	2000	75	72	4	50	39- 99
2-Nitroaniline	2778	2778	0.00	2160	1990	78	72	8	50	43-131
Dimethyl Phthalate	2778	2778	0.00	2740	2530	99	91	8	50	61-112
2,6-Dinitrotoluene	2778	2778	0.00	2640	2240	95	81	16	50	58-106
Acenaphthylene	2778	2778	0.00	2190	2040	79	73	7	50	50-115
3-Nitroaniline	2778	2778	0.00	1530	1560	55	56	2	50	25-110
Acenaphthene	2778	2778	0.00	2120	2030	76	73	4	50	50-118
2,4-Dinitrophenol	5556	5556	0.00	2830	2760	51	50	2	50	3- 93
Dibenzofuran	2778	2778	0.00	2190	2120	79	76	3	50	53-128
4-Nitrophenol	5556	5556	0.00	7060	5590	127	101	23	50	28-137
2,4-Dinitrotoluene	2778	2778	0.00	2350	2150	84	77	9	50	42-120
Diethylphthalate	2778	2778	0.00	2480	2170	89	78	13	50	57-108
Fluorene	2778	2778	0.00	2280	2100	82	76	8	50	63-105
4-Chlorophenyl-phenylether	2778	2778	0.00	2640	2500	95	90	5	50	63-110
4-Nitroaniline	2778	2778	0.00	1110	933	40	34	17	50	22- 93
4,6-Dinitro-2-Methylphenol	5556	5556	0.00	4400	4290	79	77	2	50	27-111
N-Nitrosodiphenylamine	2778	2778	0.00	2390	2190	86	79	9	50	50-115
4-Bromophenyl-phenylether	2778	2778	0.00	3500	3210	126	116	9	50	48-133
Hexachlorobenzene	2778	2778	0.00	3970	3550	143	128	11	50	33-154

KATAHDIN ANALYTICAL SERVICES
MATRIX SPIKE/MATRIX SPIKE DUPLICATE RECOVERY

Client: Woodard & Curran
Project: Old Town
PO No:
Sample Date: 05/30/06
Received Date: 05/31/06
Extraction Date: 06/04/06
Analysis Date: 06/06/06
Report Date: 06/06/2006
Matrix: SOIL

Lab ID: WG28761-3 & WG28761-4
Client ID: SS05MS & SS05MSD
SDG: WW2637
Extracted by: GN
Extraction Method: SW846 3550
Analyst: JCG
Analysis Method: SW846 8270C
Lab Prep Batch: WG28761
Units: ug/Kg

COMPOUND	MS SPIKE	MSD SPIKE	SAMPLE CONC.	MS CONC.	MSD CONC.	MS %REC.	MSD %REC.	%RPD	%RPD LIMIT	QC. LIMITS
Pentachlorophenol	5556	5556	0.00	4040	3720	73	67	8	50	32-139
Phenanthrene	2778	2778	0.00	2710	2500	98	90	8	50	55-121
Anthracene	2778	2778	0.00	2700	2570	97	92	5	50	60-122
Carbazole	2778	2778	0.00	2110	1770	76	64	18	50	51-119
Di-n-butylphthalate	2778	2778	0.00	2410	2270	87	82	6	50	40-120
Fluoranthene	2778	2778	0.00	2000	1700	72	61	16	50	41-135
Pyrene	2778	2778	0.00	2940	2670	106	96	10	50	50-114
Butylbenzylphthalate	2778	2778	0.00	2300	2000	83	72	14	50	46-115
Benzo(a)anthracene	2778	2778	0.00	2080	1980	75	71	5	50	45-111
3,3'-Dichlorobenzidine	2778	2778	0.00	1020	1140	37	41	11	50	10-130
Chrysene	2778	2778	0.00	2030	2080	73	75	2	50	45-124
bis(2-Ethylhexyl)phthalate	2778	2778	0.00	2320	2230	84	80	4	50	45-125
Di-n-octylphthalate	2778	2778	0.00	2000	1850	72	66	8	50	39-126
Benzo(b)fluoranthene	2778	2778	0.00	1400	1460	50	52	4	50	47-114
Benzo(k)fluoranthene	2778	2778	0.00	1300	1430	47	51	10	50	46-125
Benzo(a)pyrene	2778	2778	0.00	1080	1240	* 39	45	14	50	45-119
Indeno(1,2,3-cd)pyrene	2778	2778	0.00	496	646	* 18	* 23	26	50	38-131
Dibenzo(a,h)anthracene	2778	2778	0.00	582	688	* 21	* 25	17	50	37-125
Benzo(g,h,i)perylene	2778	2778	0.00	445	597	* 16	* 21	29	50	37-133

FORM III SV-2

X0018.D & X0019.D

FORM 4
PESTICIDE METHOD BLANK SUMMARY

CLIENT SAMPLE ID

| WG28624-BLANK |

Lab Name: KATAHDIN ANALYTICAL SERVICES Lab Code: KAS

Project: OLD TOWN SDG No.: WW2637

Lab Sample ID: WG28624-1 Lab File ID: 6WF1012

Matrix (soil/water) SOIL Extraction:(SepF/Cont/Sonc) SW846 3550

Sulfur Cleanup: (Y/N) N Date Extracted: 05/31/06

Date Analyzed (1): 06/01/06 Date Analyzed (2): 06/01/06

Time Analyzed (1): 1328 Time Analyzed (2): 1328

Instrument ID (1): GC06 Instrument ID (2): GC06

GC Column (1): RTX-5 ID: 0.53(mm) GC Column (2): RTX-35 ID: 0.53(mm)

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES, MS and MSD:

	CLIENT SAMPLE ID	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED 1	DATE ANALYZED 2
01	WG28624-LCS	WG28624-2	6WF1013	06/01/06	06/01/06
02	WG28624-LCSD	WG28624-3	6WF1014	06/01/06	06/01/06
03	SS07	WW2637-12	6WF1015	06/01/06	06/01/06
04	SS08	WW2637-13	6WF1016	06/01/06	06/01/06
05					
06					
07					
08					
09					
10					
11					
12					
13					
14					
15					
16					
17					
18					
19					
20					
21					
22					
23					
24					

COMMENTS: _____

KATAHDIN ANALYTICAL SERVICES
Report of Analytical Results

Client:
Project: Old Town
PO No:
Sample Date:
Received Date:
Extraction Date: 05/31/06
Analysis Date: 01-JUN-2006 13:28
Report Date: 06/02/2006
Matrix: SOIL
% Solids: 100

Lab ID: WG28624-1
Client ID: WG28624-Blank
SDG: WW2637
Extracted by: KM
Extraction Method: SW846 3550
Analyst: JLP
Analysis Method: SW846 8082
Lab Prep Batch: WG28624
Units: ug/Kg

Compound	Flags	Results	DF	PQL	Adj.PQL
Aroclor-1016	U	17	1.0	17	17
Aroclor-1221	U	17	1.0	17	17
Aroclor-1232	U	17	1.0	17	17
Aroclor-1242	U	17	1.0	17	17
Aroclor-1248	U	17	1.0	17	17
Aroclor-1254	U	17	1.0	17	17
Aroclor-1260	U	17	1.0	17	17
Tetrachloro-m-xylene		114%			
Decachlorobiphenyl		122%			

KATAHDIN ANALYTICAL SERVICES
LAB CONTROL SAMPLE

Client:
Project: Old Town
PO No:
Sample Date:
Received Date:
Extraction Date: 05/31/06
Analysis Date: 06/01/06
Report Date: 06/02/2006
Matrix: SOIL

Lab ID: WG28624-2 & WG28624-3
Client ID: WG28624-LCS & WG28624-LCSD
SDG: WW2637
Extracted by: KM
Extraction Method: SW846 3550
Analyst: JLP
Analysis Method: SW846 8082
Lab Prep Batch: WG28624
Units: ug/Kg

COMPOUND	LCS SPIKE	LCSD SPIKE	SAMPLE CONC.	LCS CONC.	LCSD CONC.	LCS %REC.		LCSD %REC.	%RPD	%RPD LIMIT	QC. LIMITS
Aroclor-1016	167	167	NA	172	185	103		111	7	50	71-114
Aroclor-1260	167	167	NA	185	202	111	*	121	9	50	77-119

FORM III PEST-2

6WF1013.d & 6WF1014.d

Katahdin
ANALYTICAL SERVICES

PREPARATION BLANK REPORT

Sample ID: PBSWE31ICS0 **Batch ID** WE31ICS0

Element Name	Result	Units	Flag	PQL	File
ALUMINUM	0.8	mg/kgdrywt	J	30.0	BWE31A
ANTIMONY	4.5	mg/kgdrywt	H	0.800	AWF01A
ARSENIC	0.5	mg/kgdrywt	J	0.800	AWF01A
BARIUM	0.1	mg/kgdrywt	U	0.500	BWE31A
BERYLLIUM	0.03	mg/kgdrywt	U	0.500	BWE31A
BORON	0.7	mg/kgdrywt	U	10.0	BWE31A
CADMIUM	0.4	mg/kgdrywt	U	1.00	BWE31A
CALCIUM	16.4	mg/kgdrywt	H	5.00	FWF02A
CHROMIUM	0.2	mg/kgdrywt	J	1.50	AWF01A
COBALT	0.3	mg/kgdrywt	U	3.00	BWE31A
COPPER	0.3	mg/kgdrywt	U	2.50	BWE31A
IRON	4.2	mg/kgdrywt	J	10.0	BWE31A
LEAD	0.2	mg/kgdrywt	U	0.500	AWF01A
MAGNESIUM	0.3	mg/kgdrywt	J	5.00	FWF02A
MANGANESE	0.25	mg/kgdrywt	J	0.500	BWE31A
MOLYBDENUM	0.2	mg/kgdrywt	U	1.00	AWF01A
NICKEL	1.	mg/kgdrywt	U	4.00	BWE31A
POTASSIUM	40.	mg/kgdrywt	U	100.	BWE31A
SELENIUM	0.4	mg/kgdrywt	U	1.00	AWF01A
SILICON	5.	mg/kgdrywt	J	20.0	BWE31A
SILVER	0.4	mg/kgdrywt	U	1.50	BWE31A
SODIUM	16.	mg/kgdrywt	J	100.	BWE31A
STRONTIUM	0.0206	mg/kgdrywt	J	10.0	FWF02A
THALLIUM	0.6	mg/kgdrywt	U	1.50	AWF01A
TIN	1.14	mg/kgdrywt	J	10.0	FWF02A
TITANIUM	0.52	mg/kgdrywt	J	1.50	AWF01A
URANIUM	0.0008	mg/kgdrywt	J	0.500	FWF02A
VANADIUM	0.5	mg/kgdrywt	U	2.50	BWE31A
ZINC	0.3	mg/kgdrywt	J	2.50	BWE31A

U The analyte was not detected in the sample at a level greater than the instrument detection limit.

J The analyte was detected in the sample at a concentration greater than the instrument detection limit, but less than the laboratory's Practical Quantitation Level.

H The analyte was detected in the sample at a concentration greater than the laboratory's acceptance limit.

Katahdin
ANALYTICAL SERVICES

LABORATORY CONTROL SAMPLE REPORT

Sample ID: LCSOWE31ICS0 **Batch ID** WE31ICS0

Element Name	True Value	Result	Units	Recovery(%)	Flag	Limits (mg/kgdrywt)		File
ALUMINUM	2.00	197.	mg/kgdrywt	98.5%		159	241	BWE31A
ANTIMONY	0.500	45.4	mg/kgdrywt	90.8%		39.8	60.2	AWF01A
ARSENIC	0.500	51.6	mg/kgdrywt	103.2%		39.8	60.2	AWF01A
BARIUM	2.00	196.	mg/kgdrywt	98.0%		159	241	BWE31A
BERYLLIUM	0.0500	4.77	mg/kgdrywt	95.4%		3.98	6.02	BWE31A
BORON	0.500	51.7	mg/kgdrywt	103.4%		39.8	60.2	BWE31A
CADMIUM	0.250	24.2	mg/kgdrywt	96.8%		19.9	30.1	BWE31A
CALCIUM	2.50	224.	mg/kgdrywt	89.6%		199	301	FWF02A
CHROMIUM	0.200	22.9	mg/kgdrywt	114.5%		15.9	24.1	AWF01A
COBALT	0.500	48.9	mg/kgdrywt	97.8%		39.8	60.2	BWE31A
COPPER	0.250	23.7	mg/kgdrywt	94.8%		199	30.1	BWE31A
IRON	1.00	103.	mg/kgdrywt	103.0%		79.5	120	BWE31A
LEAD	0.500	59.0	mg/kgdrywt	118.0%		39.8	60.2	AWF01A
MAGNESIUM	5.00	486.	mg/kgdrywt	97.2%		398	602	FWF02A
MANGANESE	0.500	49.0	mg/kgdrywt	98.0%		39.8	60.2	BWE31A
MOLYBDENUM	0.300	31.0	mg/kgdrywt	103.3%		23.8	36.1	AWF01A
NICKEL	0.50	49.	mg/kgdrywt	98.0%		39.8	60.2	BWE31A
POTASSIUM	10.0	980.	mg/kgdrywt	98.0%		795	1200	BWE31A
SELENIUM	0.500	49.0	mg/kgdrywt	98.0%		39.8	60.2	AWF01A
SILICON	0.98	142.	mg/kgdrywt	144.9%	H	416	630	BWE31A
SILVER	0.050	4.4	mg/kgdrywt	88.0%		3.98	6.02	BWE31A
SODIUM	7.50	757.	mg/kgdrywt	100.9%		596	904	BWE31A
STRONTIUM	0.750	71.3	mg/kgdrywt	95.1%		39.8	60.2	FWF02A
THALLIUM	0.500	54.0	mg/kgdrywt	108.0%		39.8	60.2	AWF01A
TIN	0.750	65.8	mg/kgdrywt	87.7%		39.8	60.2	FWF02A
TITANIUM	1.00	97.3	mg/kgdrywt	97.3%		79.5	120	AWF01A
URANIUM	0.100	9.62	mg/kgdrywt	96.2%		7.95	12.0	FWF02A
VANADIUM	0.500	48.9	mg/kgdrywt	97.8%		39.8	60.2	BWE31A
ZINC	0.500	48.1	mg/kgdrywt	96.2%		39.8	60.2	BWE31A

H Laboratory control sample recovery is greater than the laboratory's acceptance limit.

L Laboratory control sample recovery is less than the laboratory's acceptance limit.

Katahdin
ANALYTICAL SERVICES

PREPARATION BLANK REPORT

Sample ID: PBWWE31ICW1 **Batch ID** WE31ICW1

Element Name	Result	Units	Flag	PQL	File
ALUMINUM	0.04	mg/L	J	0.30	AWE31A
ANTIMONY	0.004	mg/L	U	0.008	AWE31A
ARSENIC	0.003	mg/L	U	0.008	AWF01A
BARIUM	0.0003	mg/L	U	0.0050	AWE31A
BERYLLIUM	0.0004	mg/L	U	0.0050	AWE31A
BORON	0.02	mg/L	U	0.10	AWE31A
CADMIUM	0.0004	mg/L	U	0.0100	AWE31A
CALCIUM	0.01	mg/L	U	0.05	AWE31A
CHROMIUM	0.001	mg/L	U	0.015	AWF01A
COBALT	0.001	mg/L	U	0.030	AWE31A
COPPER	0.002	mg/L	U	0.025	AWE31A
IRON	0.02	mg/L	U	0.10	AWE31A
LEAD	0.002	mg/L	U	0.005	AWF01A
MAGNESIUM	0.01	mg/L	U	0.05	AWE31A
MANGANESE	0.0002	mg/L	U	0.0050	AWE31A
MOLYBDENUM	0.002	mg/L	U	0.010	AWE31A
NICKEL	0.002	mg/L	U	0.040	AWE31A
POTASSIUM	0.4	mg/L	U	1.0	BWF01B
SELENIUM	0.004	mg/L	U	0.010	AWE31A
SILICON	0.01	mg/L	U	0.20	BWF01B
SILVER	0.001	mg/L	U	0.015	AWE31A
SODIUM	0.12	mg/L	J	1.00	BWF01B
STRONTIUM	0.00005	mg/L	U	0.100	AWE31A
THALLIUM	0.006	mg/L	U	0.015	AWF01A
TIN	0.003	mg/L	U	0.100	AWE31A
TITANIUM	0.0006	mg/L	U	0.0150	AWE31A
URANIUM	0.00001	mg/L	J	0.00500	FWF02A
VANADIUM	0.001	mg/L	U	0.025	AWE31A
ZINC	0.0012	mg/L	J	0.0250	AWE31A

U The analyte was not detected in the sample at a level greater than the instrument detection limit.

J The analyte was detected in the sample at a concentration greater than the instrument detection limit, but less than the laboratory's Practical Quantitation Level.

H The analyte was detected in the sample at a concentration greater than the laboratory's acceptance limit.

Katahdin
ANALYTICAL SERVICES

LABORATORY CONTROL SAMPLE REPORT

Sample ID: LCSWWE31ICW1 **Batch ID** WE31ICW1

Element Name	True Value	Result	Units	Recovery(%)	Flag	Limits (%)		File
ALUMINUM	2.00	2.27	mg/L	113.5%		80.	120.	AWE31A
ANTIMONY	0.500	0.477	mg/L	95.4%		80.	120.	AWE31A
ARSENIC	0.500	0.504	mg/L	100.8%		80.	120.	AWF01A
BARIUM	2.00	2.12	mg/L	106.0%		80.	120.	AWE31A
BERYLLIUM	0.0500	0.0466	mg/L	93.2%		80.	120.	AWE31A
BORON	0.50	0.45	mg/L	90.0%		80.	120.	AWE31A
CADMIUM	0.250	0.245	mg/L	98.0%		80.	120.	AWE31A
CALCIUM	2.50	2.62	mg/L	104.8%		80.	120.	AWE31A
CHROMIUM	0.200	0.215	mg/L	107.5%		80.	120.	AWF01A
COBALT	0.500	0.496	mg/L	99.2%		80.	120.	AWE31A
COPPER	0.250	0.261	mg/L	104.4%		80.	120.	AWE31A
IRON	1.00	1.06	mg/L	106.0%		80.	120.	AWE31A
LEAD	0.500	0.507	mg/L	101.4%		80.	120.	AWF01A
MAGNESIUM	5.00	5.31	mg/L	106.2%		80.	120.	AWE31A
MANGANESE	0.500	0.512	mg/L	102.4%		80.	120.	AWE31A
MOLYBDENUM	0.300	0.298	mg/L	99.3%		80.	120.	AWE31A
NICKEL	0.500	0.499	mg/L	99.8%		80.	120.	AWE31A
POTASSIUM	10.0	11.0	mg/L	110.0%		80.	120.	BWF01B
SELENIUM	0.500	0.461	mg/L	92.2%		80.	120.	AWE31A
SILICON	0.98	1.04	mg/L	106.1%		80.	120.	BWF01B
SILVER	0.050	0.048	mg/L	96.0%		80.	120.	AWE31A
SODIUM	7.50	8.15	mg/L	108.7%		80.	120.	BWF01B
STRONTIUM	0.750	0.811	mg/L	108.1%		80.	120.	AWF01A
THALLIUM	0.500	0.502	mg/L	100.4%		80.	120.	AWE31A
TIN	0.750	0.720	mg/L	96.0%		80.	120.	AWF01A
TITANIUM	1.00	1.05	mg/L	105.0%		80.	120.	AWE31A
URANIUM	0.100	0.105	mg/L	105.0%		80.	120.	FWF02A
VANADIUM	0.500	0.496	mg/L	99.2%		80.	120.	AWE31A
ZINC	0.500	0.478	mg/L	95.6%		80.	120.	AWE31A

H Laboratory control sample recovery is greater than the laboratory's acceptance limit.

L Laboratory control sample recovery is less than the laboratory's acceptance limit.

Katahdin
ANALYTICAL SERVICES

PREPARATION BLANK REPORT

Sample ID: PBSWF01HGS0 **Batch ID** WF01HGS0

Element Name	Result	Units	Flag	PQL	File
MERCURY	0.002	ug/gdrywt	U	0.04	HWF02A

U The analyte was not detected in the sample at a level greater than the instrument detection limit.

J The analyte was detected in the sample at a concentration greater than the instrument detection limit, but less than the laboratory's Practical Quantitation Level.

H The analyte was detected in the sample at a concentration greater than the laboratory's acceptance limit.

Katahdin
ANALYTICAL SERVICES

LABORATORY CONTROL SAMPLE REPORT

Sample ID: LCSOWF01HGS0 **Batch ID** WF01HGS0

Element Name	True Value	Result	Units	Recovery(%)	Flag	Limits (ug/gdrywt)		File
MERCURY	5.00	0.808	ug/gdrywt	97.0%		0.663	1.00	HWF02A

H Laboratory control sample recovery is greater than the laboratory's acceptance limit.

L Laboratory control sample recovery is less than the laboratory's acceptance limit.

Katahdin
ANALYTICAL SERVICES

PREPARATION BLANK REPORT

Sample ID: PBWWF01HGW0 **Batch ID** WF01HGW0

Element Name	Result	Units	Flag	PQL	File
MERCURY	0.02	ug/L	U	0.20	HWF01B

U The analyte was not detected in the sample at a level greater than the instrument detection limit.

J The analyte was detected in the sample at a concentration greater than the instrument detection limit, but less than the laboratory's Practical Quantitation Level.

H The analyte was detected in the sample at a concentration greater than the laboratory's acceptance limit.

Katahdin
ANALYTICAL SERVICES

LABORATORY CONTROL SAMPLE REPORT

Sample ID: LCSWWF01HGW0 **Batch ID** WF01HGW0

Element Name	True Value	Result	Units	Recovery(%)	Flag	Limits (%)		File
MERCURY	5.00	4.93	ug/L	98.6%		80.	120.	HWF01B

H Laboratory control sample recovery is greater than the laboratory's acceptance limit.

L Laboratory control sample recovery is less than the laboratory's acceptance limit.

Katahdin
ANALYTICAL SERVICES

PREPARATION BLANK REPORT

Sample ID: PBSWF06HGS0 **Batch ID** WF06HGS0

Element Name	Result	Units	Flag	PQL	File
MERCURY	0.002	ug/gdrywt	U	0.04	HWF07A

U The analyte was not detected in the sample at a level greater than the instrument detection limit.

J The analyte was detected in the sample at a concentration greater than the instrument detection limit, but less than the laboratory's Practical Quantitation Level.

H The analyte was detected in the sample at a concentration greater than the laboratory's acceptance limit.

Katahdin
ANALYTICAL SERVICES

LABORATORY CONTROL SAMPLE REPORT

Sample ID: LCSOWF06HGS0 **Batch ID** WF06HGS0

Element Name	True Value	Result	Units	Recovery(%)	Flag	Limits (ug/gdrywt)		File
MERCURY	5.00	0.750	ug/gdrywt	90.0%		0.663	1.00	HWF07A

H Laboratory control sample recovery is greater than the laboratory's acceptance limit.

L Laboratory control sample recovery is less than the laboratory's acceptance limit.

Katahdin
ANALYTICAL SERVICES

MATRIX SPIKE / MATRIX SPIKE DUPLICATE QC SUMMARY

Sample ID: WW2637-001

Symbol	Sample Result		Units	Spike Added	Spike Result	Spike Rec.(%)	Note	Spike Duplicate Result	Spike Duplicate Rec.(%)	Note	RPD(%)	Note
Ag	U	0.015	mg/L	0.0500	0.048	96.8 %		0.048	96.8 %		0.00 %	
Ba		0.0774	mg/L	2.00	2.15	104. %		2.09	101. %		2.83 %	
Cd	U	0.0100	mg/L	0.250	0.260	104. %		0.249	99.5 %		4.32 %	
Cr	U	0.0030	mg/L	0.200	0.238	118. %		0.223	110. %		6.51 %	
HG	U	0.20	ug/L	1.00	0.95	99.0 %		0.96	100. %		1.05 %	
Pb	U	0.0010	mg/L	0.500	0.523	104. %		0.503	100. %		3.90 %	
Se	U	0.010	mg/L	0.500	0.515	103. %		0.489	97.4 %		5.18 %	

Matrix spike recovery is outside the laboratory's specified acceptance range indicating potential sample matrix interference and potential bias of reported value for this parameter.

Matrix spike recovery is outside the laboratory's specified acceptance range. The spike concentration for this parameter is significantly below the sample concentration and cannot be distinguished from the sample's analytical signal.

Matrix spike analysis cannot be quantified due to severe matrix interferences.

Precision of replicate analysis as measured by RPD is outside the laboratory's acceptance range for this parameter. Sample homogeneity may be a factor.

Because of the large uncertainty associated with measurements made near the detection level, there is no acceptance range for relative percent difference.

Katahdin
ANALYTICAL SERVICES

MATRIX SPIKE / MATRIX SPIKE DUPLICATE QC SUMMARY

Sample ID: WW2637-011

Symbol	Sample Result		Units	Spike Added	Spike Result	Spike Rec.(%)	Note	Spike Duplicate Result	Spike Duplicate Rec.(%)	Note	RPD(%)	Note
Ag	U	2.2	mg/Kg	6.89	6.2	98.7 %		6.7	106. %		7.75 %	
As		1.	mg/Kg	68.9	73.3	105. %		78.9	113. %		7.36 %	
Ba		85.6	mg/Kg	276.	307.	80.4 %		335.	90.5 %		8.72 %	
Cd		0.393	mg/Kg	34.4	37.7	108. %		40.2	116. %		6.42 %	
Cr		29.8	mg/Kg	27.6	57.2	99.4 %		60.4	111. %		5.44 %	
HG	U	0.058	ug/g	0.278	0.25	91.8 %		0.23	84.6 %		8.33 %	
Pb		4.43	mg/Kg	68.9	81.0	111. %		85.7	118. %		5.64 %	
Se	U	1.5	mg/Kg	68.9	70.0	103. %		74.4	109. %		6.09 %	

Matrix spike recovery is outside the laboratory's specified acceptance range indicating potential sample matrix interference and potential bias of reported value for this parameter.

Matrix spike recovery is outside the laboratory's specified acceptance range. The spike concentration for this parameter is significantly below the sample concentration and cannot be distinguished from the sample's analytical signal.

Matrix spike analysis cannot be quantified due to severe matrix interferences.

Precision of replicate analysis as measured by RPD is outside the laboratory's acceptance range for this parameter. Sample homogeneity may be a factor.

Because of the large uncertainty associated with measurements made near the detection level, there is no acceptance range for relative percent difference.

Katahdin Analytical Services WW2637 Page 0000113 of 0000126

Report of Analytical Results

Client: Erika Lloyd
Woodard & Curran
41 Hutchins Drive
Portland,ME 04102

Lab Sample ID: WW2637-6
Report Date: 07-JUN-06
Client PO:
Project: Old Town
SDG: WW2637

Sample Description	Matrix	Date Sampled	Date Received
SS01	SL	30-MAY-06	31-MAY-06

Parameter	Result	Adj PQL	Anal. Method	QC.Batch	Anal. Date	Prep. Method	Prep. Date	Analyst	Footnotes
Total Solids	45. %	1	CLP SOW 788	WG28725	02-JUN-06 08:41:00	CLP SOW 788	01-JUN-06	JF	

Report of Analytical Results

Client: Erika Lloyd
Woodard & Curran
41 Hutchins Drive
Portland,ME 04102

Lab Sample ID: WW2637-7
Report Date: 07-JUN-06
Client PO:
Project: Old Town
SDG: WW2637

Sample Description	Matrix	Date Sampled	Date Received
SS02	SL	30-MAY-06	31-MAY-06

Parameter	Result	Adj PQL	Anal. Method	QC.Batch	Anal. Date	Prep. Method	Prep. Date	Analyst	Footnotes
Total Solids	92. %	1	CLP SOW 788	WG28725	02-JUN-06 08:43:00	CLP SOW 788	01-JUN-06	JF	

Report of Analytical Results

Client: Erika Lloyd
Woodard & Curran
41 Hutchins Drive
Portland,ME 04102

Lab Sample ID: WW2637-8
Report Date: 07-JUN-06
Client PO:
Project: Old Town
SDG: WW2637

Sample Description	Matrix	Date Sampled	Date Received
SS03	SL	30-MAY-06	31-MAY-06

Parameter	Result	Adj PQL	Anal. Method	QC.Batch	Anal. Date	Prep. Method	Prep. Date	Analyst	Footnotes
Total Solids	53. %	1	CLP SOW 788	WG28725	02-JUN-06 08:44:00	CLP SOW 788	01-JUN-06	JF	

Report of Analytical Results

Client: Erika Lloyd
Woodard & Curran
41 Hutchins Drive
Portland,ME 04102

Lab Sample ID: WW2637-9
Report Date: 07-JUN-06
Client PO:
Project: Old Town
SDG: WW2637

Sample Description	Matrix	Date Sampled	Date Received
SS80	SL	30-MAY-06	31-MAY-06

Parameter	Result	Adj PQL	Anal. Method	QC.Batch	Anal. Date	Prep. Method	Prep. Date	Analyst	Footnotes
Total Solids	50. %	1	CLP SOW 788	WG28725	02-JUN-06 08:45:00	CLP SOW 788	01-JUN-06	JF	

Quality Control Report

Blank Sample Summary Report

Total Solids

Samp Type	QC Batch	Anal. Method	Anal. Date	Prep. Date	Result	PQL
MBLANK	WG28725	CLP SOW 788	02-JUN-06	01-JUN-06	U 1 %	1 %

Quality Control Report

Duplicate Sample Summary Report

Total Solids

Duplicate Sample ID	Original Sample ID	QC Batch	Analysis Date	Result Units	Sample Result	Duplicate Result	RPD(%)	RPD Limit
WG28725-3	WW2637-6	WG28725	02-JUN-06	%	45.	45.	1	20
WG28725-4	WW2637-11	WG28725	02-JUN-06	%	60.	62.	3	20

Quality Control Report

Laboratory Control Sample Summary Report

Total Solids

Lab Sample Id	Samp Type	QC Batch	Analysis Date	Prep Date	Units	Spike Amt.	Result	Recovery	Acceptance Range	RPD
WG28725-2	LCS	WG28725	02-JUN-06	01-JUN-06	%	90	90.	100	80-120	

KATAHDIN ANALYTICAL SERVICES, INC.
SAMPLE RECEIPT CONDITION REPORT
Tel. (207) 874-2400
Fax (207) 775-4029

PAGE: 1 OF 2

COOLER: 1 OF 2

CLIENT: W&L

PROJECT:

COC#

SDG#

DATE / TIME RECEIVED: 5-31-06 0900

DELIVERED BY: Client

RECEIVED BY: AA.

LIMS ENTRY BY: DWM

LIMS REVIEW BY / PM: AJC

	YES	NO	EXCEPTIONS
1. CUSTODY SEALS PRESENT / INTACT?	☒	☐	☐
2. CHAIN OF CUSTODY PRESENT IN THIS COOLER?	☒	☐	☐
3. CHAIN OF CUSTODY SIGNED BY CLIENT?	☒	☐	☐
4. CHAIN OF CUSTODY MATCHES SAMPLES?	☐	☒	☐
5. TEMPERATURE BLANKS PRESENT?	☒	☐	☐
6. SAMPLES RECEIVED AT 4°C +/- 2?	☒	☐	☐
ICE ICE PACKS PRESENT (Y) or N?			
7. VOLATILES FREE OF HEADSPACE?	☒	☐	☐
8. TRIP BLANK PRESENT IN THIS COOLER	☒	☐	☐
9. PROPER SAMPLE CONTAINERS AND VOLUME?	☒	☐	☐
10. SAMPLES WITHIN HOLD TIME UPON RECEIPT?	☒	☐	☐
11. SAMPLES PROPERLY PRESERVED[1]?	☒	☐	☐
12. CORRECTIVE ACTION REPORT FILED?	☐	☒	N/A

COMMENTS / RESOLUTION

TEMP BLANK TEMP (°C)= 2.3

COOLER TEMP (°C)= NA
(RECORD COOLER TEMP ONLY IF TEMP BLANK IS NOT PRESENT)

13. ANALYTICAL PROGRAMS (CIRCLE ONE) (COMMERCIAL) CLP HAZWRAP NFESC ACOE AFCEE OTHER (STATE OF ORIGIN):

LOG - IN NOTES[1]: Sample 5581 not on COC, logged in as WW2637-14 - Runs on Stone Bartlett 5/31/06.
VOA Trip Blank not on COC, logged in as WW2637-15.

[1] Use this space (and additional sheets if necessary) to document samples that are received broken or compromised, C-O-C discrepancies, radiation checks, residual chlorine check, results of pH check if required. If samples required pH adjustment, record volume and type of preservative added

Katahdin Analytical Services WW2637 Page 0000121 of 0000126

KATAHDIN ANALYTICAL SERVICES, INC.

SAMPLE RECEIPT CONDITION REPORT

Tel. (207) 874-2400
Fax (207) 775-4029

PAGE: 2 OF 2

COOLER: 2 OF 2

CLIENT: W&L

PROJECT:

COC#:
SDG#:
DATE / TIME RECEIVED: 5-31-06 0900
DELIVERED BY: Client
RECEIVED BY: AA
LIMS ENTRY BY: DWM
LIMS REVIEW BY / PM: AJC

	YES	NO	EXCEPTIONS
1. CUSTODY SEALS PRESENT / INTACT?	☒	☐	☐
2. CHAIN OF CUSTODY PRESENT IN THIS COOLER?	☒	☐	☐
3. CHAIN OF CUSTODY SIGNED BY CLIENT?	☒	☐	☐
4. CHAIN OF CUSTODY MATCHES SAMPLES?	☐	☒	☐
5. TEMPERATURE BLANKS PRESENT?	☒	☐	☐
6. SAMPLES RECEIVED AT 4°C +/- 2?	☒	☐	☐
7. ICE / ICE PACKS PRESENT (Y or N)?			
8. VOLATILES FREE OF HEADSPACE?	☐	☐	☒
9. TRIP BLANK PRESENT IN THIS COOLER	☐	☐	☒
10. PROPER SAMPLE CONTAINERS AND VOLUME?	☒	☐	☐
11. SAMPLES WITHIN HOLD TIME UPON RECEIPT?	☒	☐	☐
12. SAMPLES PROPERLY PRESERVED[1]?	☒	☐	☐
13. CORRECTIVE ACTION REPORT FILED?	☐	☒	N/A

COMMENTS / RESOLUTION

TEMP BLANK TEMP (°C)= 7.4

COOLER TEMP (°C)= 6.0 NA

(RECORD COOLER TEMP ONLY IF TEMP BLANK IS NOT PRESENT)

14. ANALYTICAL PROGRAMS (CIRCLE ONE) COMMERCIAL CLP HAZWRAP NFESC ACOE AFCEE OTHER (STATE OF ORIGIN):

LOG - IN NOTES[1]:

[1] Use this space (and additional sheets if necessary) to document samples that are received broken or compromised, C-O-C discrepancies, radiation checks, residual chlorine check, results of pH check if required. If samples required pH adjustment, record volume and type of preservative added

Katahdin Analytical Services WW2637 Page 0000122 of 0000126

Katahdin ANALYTICAL SERVICES
340 County Road No. 5
P.O. Box 720
Westbrook, ME 04092
Tel: (207) 874-2400
Fax: (207) 775-4029

CHAIN of CUSTODY

PLEASE BEAR DOWN AND PRINT LEGIBLY IN PEN

Page ____ of ____

Client	Contact	Phone #	Fax #
Woodard + Curran	KARI KASPER	(207) 774-2112	(207) 774-6635

Address	City	State	Zip Code
41 Hutchins DR	Portland	ME	04102

Purchase Order # | Proj. Name / No. | Katahdin Quote #

Bill (if different than above) | Address

Sampler (Print / Sign) DANIELLE TAYLOR Danielle Taylor | Copies To:

LAB USE ONLY WORK ORDER #: WW2637
KATAHDIN PROJECT NUMBER ____

REMARKS: ____

SHIPPING INFO: ☐ FED EX ☐ UPS ☐ CLIENT

AIRBILL NO: ____

TEMP°C ____ ☐ TEMP BLANK ☐ INTACT ☐ NOT INTACT

*	Sample Description	Date / Time coll'd	Matrix	No. of Cntrs.	ANALYSIS AND CONTAINER TYPE PRESERVATIVES: Filt. ☐Y ☑N — SVOCS 8270	Filt. ☑Y ☐N — RCRA 8 Metals	Filt. ☐Y ☑N — VOCs 8260B	Filt. ☑Y ☐N — PCBs	Filt. ☐Y ☐N	Filt. ☐Y ☐N	Filt. ☐Y ☐N	Filt. ☐Y ☐N	Filt. ☐Y ☐N	Filt. ☐Y ☐N
	SB-25	5/30/06/13:55	GW	6	✓	✓	✓							
	MW-1	/0855	GW	6	✓	✓	✓							
	MW-2	/11:10	GW	6	✓	✓	✓							
	SB-25 MS	/13:55	GW	6	✓	✓	✓							
	SB-25 MSD	/13:55	GW	6	✓	✓	✓							
	SB-26	/1450	GW	6	✓	✓	✓							
	SB-100	/1400	GW	6	✓	✓	✓							
	SS01	/0852	S	6	✓	✓	✓							
	SS02	/10:15	S	1	✓									
	SS03	/10:50	S	6	✓	✓	✓							
	SS80	/10:50	S	2	✓	✓								
	SS04	/11:30	S	6	✓	✓	✓							
	SS05	/1200	S	6	✓	✓	✓							
	SS05 MS	/1200	S	6	✓	✓	✓							
	SS07	/1230	S	1				✓						
	SS08	/1230	S	1				✓						
	SS05 MSD	1200	S	6	✓	✓	✓							

COMMENTS 48 HR TAT (Results due 7th)

Relinquished By: (Signature)	Date / Time	Received By: (Signature)	Relinquished By: (Signature)	Date / Time	Received By: (Signature)
Danielle Taylor	5/31/06 0830	AMY Adams			
Relinquished By: (Signature)	Date / Time	Received By: (Signature)	Relinquished By: (Signature)	Date / Time	Received By: (Signature)

THE TERMS AND CONDITIONS ON THE REVERSE SIDE HEREOF SHALL GOVERN

Katahdin
ANALYTICAL SERVICES

Katahdin Analytical Services

Login Chain of Custody Report (lno1)

Jun. 01, 2006
05:57 PM

Login Number: WW2637

Account: WOODAR001 Web
Woodard & Curran

Project:

Primary Report Address:
Erika Lloyd
Woodard & Curran
41 Hutchins Drive

Portland,ME 04102

Primary Invoice Address:
Accounts Payable
Woodard & Curran
41 Hutchins Drive

Portland,ME 04102

Report CC Addresses:
Invoice CC Addresses:

Login Information

ANALYSIS INSTRUCTIONS	:
CHECK NO.	:
CLIENT PO#	:
COOLER TEMPERATURE	: 2.3, 6.0
DELIVERY SERVICES	: Client
EDD FORMAT	: KAS064-XLS & KAS027-XLS
MAIL DATE	:
PM	: AJC
PROJECT NAME	: Old Town
QC LEVEL	: II
REGULATORY LIST	:
REPORT INSTRUCTIONS	: Merge results for EDDs.
SDG ID	:
SDG STATUS	:

Laboratory Sample ID	Client Sample Number	Collect Date/Time	Receive Date	PR	Verbal Date	Due Date	Comments
WW2637-1	SB-25	30-MAY-06 13:55	31-MAY-06		**02-JUN-06**	07-JUN-06	MS/MSD, limited volume for SVOA analysis

Matrix		Product	Hold Date (shortest)	Bottle Type	Bottle Count
Aqueous	P	RCRA-METALS			3
		SW3010-PREP, SW6010-ARSENIC, SW6010-BARIUM, SW6010-CADMIUM, SW6010-CHROMIUM, SW6010-LEAD, SW6010-SELENIUM, SW6010-SILVER, SW7470-MERCURY			
Aqueous	S	SW8260FULL5ML	13-JUN-06	40mL Vial+HCl	9
Aqueous	S	SW8270BNA	06-JUN-06	1L N-Amber Glass	6

Laboratory Sample ID	Client Sample Number	Collect Date/Time	Receive Date	PR	Verbal Date	Due Date	Comments
WW2637-2	MW-1	30-MAY-06 08:55	31-MAY-06		**02-JUN-06**	07-JUN-06	Limited volume for SVOA analysis

Matrix		Product	Hold Date (shortest)	Bottle Type	Bottle Count
Aqueous	P	RCRA-METALS			1
		SW3010-PREP, SW6010-ARSENIC, SW6010-BARIUM, SW6010-CADMIUM, SW6010-CHROMIUM, SW6010-LEAD, SW6010-SELENIUM, SW6010-SILVER, SW7470-MERCURY			
Aqueous	S	SW8260FULL5ML	13-JUN-06	40mL Vial+HCl	3
Aqueous	S	SW8270BNA	06-JUN-06	1L N-Amber Glass	2

Laboratory Sample ID	Client Sample Number	Collect Date/Time	Receive Date	PR	Verbal Date	Due Date	Comments
WW2637-3	MW-2	30-MAY-06 11:10	31-MAY-06		**02-JUN-06**	07-JUN-06	Limited volume for SVOA analysis

Matrix		Product	Hold Date (shortest)	Bottle Type	Bottle Count
Aqueous	P	RCRA-METALS			1
		SW3010-PREP, SW6010-ARSENIC, SW6010-BARIUM, SW6010-CADMIUM, SW6010-CHROMIUM, SW6010-LEAD, SW6010-SELENIUM, SW6010-SILVER, SW7470-MERCURY			
Aqueous	S	SW8260FULL5ML	13-JUN-06	40mL Vial+HCl	3
Aqueous	S	SW8270BNA	06-JUN-06	1L N-Amber Glass	2

Laboratory Sample ID	Client Sample Number	Collect Date/Time	Receive Date	PR	Verbal Date	Due Date	Comments
WW2637-4	SB-26	30-MAY-06 14:50	31-MAY-06		**02-JUN-06**	07-JUN-06	Limited volume for SVOA analysis

Matrix		Product	Hold Date (shortest)	Bottle Type	Bottle Count
Aqueous	P	RCRA-METALS			1
		SW3010-PREP, SW6010-ARSENIC, SW6010-BARIUM, SW6010-CADMIUM, SW6010-CHROMIUM, SW6010-LEAD, SW6010-SELENIUM, SW6010-SILVER, SW7470-MERCURY			
Aqueous	S	SW8260FULL5ML	13-JUN-06	40mL Vial+HCl	3
Aqueous	S	SW8270BNA	06-JUN-06	1L N-Amber Glass	2

Login Number: WW2637

Account: WOODAR001 Web

Woodard & Curran

Project:

Laboratory Sample ID	Client Sample Number	Collect Date/Time	Receive Date	PR	Verbal Date	Due Date	Comments
WW2637-5	SB-100	30-MAY-06 14:00	31-MAY-06		**02-JUN-06**	07-JUN-06	Limited volume for SVOA analysis

Matrix		Product	Hold Date (shortest)	Bottle Type	Bottle Count
Aqueous	P	RCRA-METALS			1
		SW3010-PREP SW6010-ARSENIC SW6010-BARIUM SW6010-CADMIUM SW6010-CHROMIUM SW6010-LEAD SW6010-SELENIUM SW6010-SILVER SW7470-MERCURY			
Aqueous	S	SW8260FULL5ML	13-JUN-06	40mL Vial+HCl	3
Aqueous	S	SW8270BNA	06-JUN-06	1L N-Amber Glass	2
WW2637-6		SS01	30-MAY-06 08:52	31-MAY-06	**02-JUN-06**	07-JUN-06	
Solid	P	RCRA-METALS			1
		SW3050-PREP SW6010-ARSENIC SW6010-BARIUM SW6010-CADMIUM SW6010-CHROMIUM SW6010-LEAD SW6010-SELENIUM SW6010-SILVER SW7471-MERCURY			
Solid	S	SW8260FULL5ML	13-JUN-06	40mL Vial+HCl	4
Solid	S	SW8270BNA	13-JUN-06	1L N-Amber Glass	1
Solid	S	TS	29-JUN-06		
WW2637-7		SS02	30-MAY-06 10:15	31-MAY-06	**02-JUN-06**	07-JUN-06	
Solid	S	SW8270BNA	13-JUN-06	1L N-Amber Glass	1
Solid	S	TS	29-JUN-06		
WW2637-8		SS03	30-MAY-06 10:50	31-MAY-06	**02-JUN-06**	07-JUN-06	
Solid	P	RCRA-METALS			1
		SW3050-PREP SW6010-ARSENIC SW6010-BARIUM SW6010-CADMIUM SW6010-CHROMIUM SW6010-LEAD SW6010-SELENIUM SW6010-SILVER SW7471-MERCURY			
Solid	S	SW8260FULL5ML	13-JUN-06	40mL Vial+HCl	4
Solid	S	SW8270BNA	13-JUN-06	1L N-Amber Glass	1
Solid	S	TS	29-JUN-06		
WW2637-9		SS80	30-MAY-06 10:50	31-MAY-06	**02-JUN-06**	07-JUN-06	
Solid	P	RCRA-METALS			1
		SW3050-PREP SW6010-ARSENIC SW6010-BARIUM SW6010-CADMIUM SW6010-CHROMIUM SW6010-LEAD SW6010-SELENIUM SW6010-SILVER SW7471-MERCURY			
Solid	S	SW8270BNA	13-JUN-06	1L N-Amber Glass	1
Solid	S	TS	29-JUN-06		
WW2637-10		SS04	30-MAY-06 11:30	31-MAY-06	**02-JUN-06**	07-JUN-06	
Solid	P	RCRA-METALS			1
		SW3050-PREP SW6010-ARSENIC SW6010-BARIUM SW6010-CADMIUM SW6010-CHROMIUM SW6010-LEAD SW6010-SELENIUM SW6010-SILVER SW7471-MERCURY			
Solid	S	SW8260FULL5ML	13-JUN-06	40mL Vial+HCl	4
Solid	S	SW8270BNA	13-JUN-06	1L N-Amber Glass	1
Solid	S	TS	29-JUN-06		

Login Number: WW2637

Account: WOODAR001 Web
Woodard & Curran

Project:

Laboratory Sample ID	Client Sample Number	Collect Date/Time	Receive Date	PR	Verbal Date	Due Date	Comments
WW2637-11	SS05	30-MAY-06 12:00	31-MAY-06		**02-JUN-06**	07-JUN-06	MS/MSD
Matrix	*Product*	*Hold Date (shortest)*	*Bottle Type*		*Bottle Count*		
Solid	P RCRA-METALS				3		
	SW3050-PREP SW6010-ARSENIC SW6010-BARIUM SW6010-CADMIUM SW6010-CHROMIUM SW6010-LEAD SW6010-SELENIUM SW6010-SILVER SW7471-MERCURY						
Solid	S SW8260FULL5ML	13-JUN-06	40mL Vial+HCl		12		
Solid	S SW8270BNA	13-JUN-06	1L N-Amber Glass		3		
Solid	S TS	29-JUN-06					
WW2637-12	SS07	30-MAY-06 12:30	31-MAY-06		**02-JUN-06**	07-JUN-06	
Matrix	*Product*	*Hold Date (shortest)*	*Bottle Type*		*Bottle Count*		
Solid	S SW8082	13-JUN-06	100g Glass		1		
Solid	S TS	29-JUN-06					
WW2637-13	SS08	30-MAY-06 12:30	31-MAY-06		**02-JUN-06**	07-JUN-06	
Matrix	*Product*	*Hold Date (shortest)*	*Bottle Type*		*Bottle Count*		
Solid	S SW8082	13-JUN-06	100g Glass		1		
Solid	S TS	29-JUN-06					
WW2637-14	SS81	30-MAY-06 12:00	31-MAY-06		**02-JUN-06**	07-JUN-06	No TS aliquot, rpt on wet wt basis.
Matrix	*Product*	*Hold Date (shortest)*	*Bottle Type*		*Bottle Count*		
Solid	S SW8260FULL5ML	13-JUN-06	40mL Vial+HCl		4		
WW2637-15	TRIP BLANK	09-MAY-06 00:00	31-MAY-06		**02-JUN-06**	07-JUN-06	
Matrix	*Product*	*Hold Date (shortest)*	*Bottle Type*		*Bottle Count*		
Aqueous	S SW8260FULL5ML	23-MAY-06	40mL Vial+HCl		2		

Total Samples: **15** ***Total Analyses:*** **42**

Katahdin
ANALYTICAL SERVICES

nelac
Cert. No. E87604

June 8, 2006

Ms. Erika Lloyd
Woodard & Curran
41 Hutchins Drive
Portland,ME 04102

RE: Katahdin Lab Number: WW2676
Project ID: 218736
Project Manager: Mrs. Andrea Colby
Sample Receipt Date(s): June 01, 2006

Dear Ms. Lloyd:

Please find enclosed the following information:

* Report of Analysis (Analytical and/or Field)
* Quality Control Data Summary
* Chain of Custody (COC)
* Login Report

A copy of the Chain of Custody is included in the paginated report. The original COC is attached as an addendum to this report.

Should you have any questions or comments concerning this Report of Analysis, please do not hesitate to contact the project manager listed above. This cover letter is an integral part of the ROA.

We certify that the test results provided in this report meet all the requirements of the NELAC standards unless otherwise noted in an attached technical narrative or in the Report of Analysis.

We appreciate your continued use of our laboratory and look forward to working with you in the future. The following signature indicates technical review and acceptance of the data.

Sincerely,

KATAHDIN ANALYTICAL SERVICES

Jessica Dimond
Authorized Signature

6.9.06
Date

TECHNICAL NARRATIVE

Organics Laboratory

The sample of work order WW2676 was analyzed in accordance with "Test Methods for Evaluating Solid Wastes: Physical/Chemical Methods." SW-846. 2nd edition, 1982 (revised 1984), 3rd edition, 1986, and Updates I, II, IIA, and III 1996, Office of Solid Waste and Emergency Response, U.S. EPA, for the specific methods listed below or on the Report of Analysis. Some manual integrations may have been performed due to split peaks and/or corrected baselines. All have been flagged with an "M" (software-generated) on the pertinent quantitation reports.

8270C Analysis

The reported percent recovery acceptance limits for the Laboratory Control Samples (LCSs) are statistically derived for the full list of spiked compounds. The recoveries of the spiked analytes in the LCS, Matrix Spike (MS) and Matrix Spike Duplicate (MSD) are compared to these acceptance limits. Katahdin standard operating procedure is not to take corrective action until greater than ten percent of the spiked analytes (with the exception of the common laboratory phthalate ester contaminants) in the LCS are outside of the QC limits. If the associated MS/MSD have greater than ten percent of the spiked analytes outside of the QC limits, no corrective action is taken, as long the LCS is acceptable.

Sample WW2676-1 had a low response for one internal standard, which was outside the laboratory acceptance limit of -50% to 100% of the response of the internal standard of the daily continuing calibration standard. Based on the sample chromatogram, this deviation is likely attributable to a matrix effect. Therefore, the sample was not reanalyzed.

The laboratory method blank WG28762-1 had a high concentration for the analyte bis(2-ethylhexyl)phthalate. Since this analyte was not detected in the sample, no further action was taken.

There were no other protocol deviations or observations noted by the organics laboratory staff.

DATA QUALIFIERS

U	Indicates the compound was analyzed for but not detected above the laboratory Practical Quantitation Limit.
*	Compound recovery outside of quality control limits.
D	Indicates the result was obtained from analysis of a diluted sample. Surrogate recoveries may not be calculable.
E	Estimated value. This flag identifies compounds whose concentrations exceed the upper level of the calibration range of the instrument for that specific analysis.
J	Estimated value. The analyte was detected in the sample at a concentration less than the laboratory Practical Quantitation Limit (PQL), but above the Method Detection Limit (MDL).
B	Organics- Indicates the analyte was detected in the laboratory method blank analyzed concurrently with the sample. Metals- Indicates the analyte was detected in the sample at a concentration greater than the instrument detection limit, but less than the laboratory's Practical Quantitation Level.
N	Presumptive evidence of a compound based on a mass spectral library search.
A	Indicates that a tentatively identified compound is a suspected aldol-condensation product.
P	Used for Pesticide/Aroclor analyte when there is a greater than 25% difference for detected concentrations between the two GC columns.
MCL	Maximum Contaminant Level
NL	No limit
NFL	No Free Liquid Present
FLP	Free Liquid Present

KATAHDIN ANALYTICAL SERVICES
Report of Analytical Results

Client: Woodard & Curran
Project: 218736
PO No:
Sample Date: 06/01/06
Received Date: 06/01/06
Extraction Date: 06/04/06
Analysis Date: 06-JUN-2006 20:21
Report Date: 06/07/2006
Matrix: WATER
% Solids: NA

Lab ID: WW2676-1
Client ID: EB-101
SDG: WW2676
Extracted by: KM
Extraction Method: SW846 3510
Analyst: JCG
Analysis Method: SW846 8270C
Lab Prep Batch: WG28762
Units: ug/L

Compound	Flags	Results	DF	PQL	Adj.PQL
Phenol	U	10	1.0	10	10
Bis(2-Chloroethyl)ether	U	10	1.0	10	10
2-Chlorophenol	U	10	1.0	10	10
1,3-Dichlorobenzene	U	10	1.0	10	10
1,4-Dichlorobenzene	U	10	1.0	10	10
1,2-Dichlorobenzene	U	10	1.0	10	10
2-Methylphenol	U	10	1.0	10	10
2,2'-Oxybis(1-Chloropropane)	U	10	1.0	10	10
N-Nitroso-di-n-propylamine	U	10	1.0	10	10
3&4-Methylphenol	U	10	1.0	10	10
Hexachloroethane	U	10	1.0	10	10
Nitrobenzene	U	10	1.0	10	10
Isophorone	U	10	1.0	10	10
2-Nitrophenol	U	10	1.0	10	10
2,4-Dimethylphenol	U	10	1.0	10	10
Bis(2-Chloroethoxy)methane	U	10	1.0	10	10
2,4-Dichlorophenol	U	10	1.0	10	10
1,2,4-Trichlorobenzene	U	10	1.0	10	10
Naphthalene	U	10	1.0	10	10
4-Chloroaniline	U	10	1.0	10	10
Hexachlorobutadiene	U	10	1.0	10	10
4-Chloro-3-Methylphenol	U	10	1.0	10	10
2-Methylnaphthalene	U	10	1.0	10	10
Hexachlorocyclopentadiene	U	10	1.0	10	10
2,4,6-Trichlorophenol	U	10	1.0	10	10
2,4,5-Trichlorophenol	U	25	1.0	25	25
2-Chloronaphthalene	U	10	1.0	10	10
2-Nitroaniline	U	25	1.0	25	25
Dimethyl Phthalate	U	10	1.0	10	10
2,6-Dinitrotoluene	U	10	1.0	10	10
Acenaphthylene	U	10	1.0	10	10
3-Nitroaniline	U	25	1.0	25	25
Acenaphthene	U	10	1.0	10	10
2,4-Dinitrophenol	U	25	1.0	25	25
Dibenzofuran	U	10	1.0	10	10
4-Nitrophenol	U	25	1.0	25	25
2,4-Dinitrotoluene	U	10	1.0	10	10
Diethylphthalate	U	10	1.0	10	10
Fluorene	U	10	1.0	10	10
4-Chlorophenyl-phenylether	U	10	1.0	10	10
4-Nitroaniline	U	25	1.0	25	25
4,6-Dinitro-2-Methylphenol	U	25	1.0	25	25
N-Nitrosodiphenylamine	U	10	1.0	10	10

KATAHDIN ANALYTICAL SERVICES
Report of Analytical Results

Client: Woodard & Curran
Project: 218736
PO No:
Sample Date: 06/01/06
Received Date: 06/01/06
Extraction Date: 06/04/06
Analysis Date: 06-JUN-2006 20:21
Report Date: 06/07/2006
Matrix: WATER
% Solids: NA

Lab ID: WW2676-1
Client ID: EB-101
SDG: WW2676
Extracted by: KM
Extraction Method: SW846 3510
Analyst: JCG
Analysis Method: SW846 8270C
Lab Prep Batch: WG28762
Units: ug/L

Compound	Flags	Results	DF	PQL	Adj.PQL
4-Bromophenyl-phenylether	U	10	1.0	10	10
Hexachlorobenzene	U	10	1.0	10	10
Pentachlorophenol	U	25	1.0	25	25
Phenanthrene	U	10	1.0	10	10
Anthracene	U	10	1.0	10	10
Carbazole	U	10	1.0	10	10
Di-n-butylphthalate	U	10	1.0	10	10
Fluoranthene	U	10	1.0	10	10
Pyrene	U	10	1.0	10	10
Butylbenzylphthalate	U	10	1.0	10	10
Benzo(a)anthracene	U	10	1.0	10	10
3,3'-Dichlorobenzidine	U	10	1.0	10	10
Chrysene	U	10	1.0	10	10
bis(2-Ethylhexyl)phthalate	U	10	1.0	10	10
Di-n-octylphthalate	U	10	1.0	10	10
Benzo(b)fluoranthene	U	10	1.0	10	10
Benzo(k)fluoranthene	U	10	1.0	10	10
Benzo(a)pyrene	U	10	1.0	10	10
Indeno(1,2,3-cd)pyrene	U	10	1.0	10	10
Dibenzo(a,h)anthracene	U	10	1.0	10	10
Benzo(g,h,i)perylene	U	10	1.0	10	10
2-Fluorophenol		34%			
Phenol-D6		37%			
Nitrobenzene-D5		79%			
2-Fluorobiphenyl		95%			
2,4,6-Tribromophenol		128%			
Terphenyl-D14		110%			

KATAHDIN ANALYTICAL SERVICES
Report of Analytical Results

Client: Woodard & Curran
Project: 218736
PO No:
Sample Date: 06/01/06
Received Date: 06/01/06
Extraction Date: 06/05/06
Analysis Date: 07-JUN-2006 02:12
Report Date: 06/07/2006
Matrix: WATER
% Solids: NA

Lab ID: WW2676-1
Client ID: EB-101
SDG: WW2676
Extracted by: KM
Extraction Method: SW846 3510
Analyst: JLP
Analysis Method: SW846 8082
Lab Prep Batch: WG28773
Units: ug/L

Compound	Flags	Results	DF	PQL	Adj.PQL
Aroclor-1016	U	0.50	1.0	0.50	0.50
Aroclor-1221	U	0.50	1.0	0.50	0.50
Aroclor-1232	U	0.50	1.0	0.50	0.50
Aroclor-1242	U	0.50	1.0	0.50	0.50
Aroclor-1248	U	0.50	1.0	0.50	0.50
Aroclor-1254	U	0.50	1.0	0.50	0.50
Aroclor-1260	U	0.50	1.0	0.50	0.50
Tetrachloro-m-xylene		81%			
Decachlorobiphenyl		60%			

Katahdin
ANALYTICAL SERVICES

REPORT OF ANALYTICAL RESULTS

Client: Erika Lloyd
Woodard & Curran
41 Hutchins Drive
Portland, ME 04102

Lab Sample ID: WW2676-001
Report Date: 6/8/2006
PO No.:
Project: 218736

Sample Description	Matrix	Filtered	Date Sampled	Date Received
EB-101	AQ	No(Total)	06/01/2006	06/01/2006

Parameter	Result	Units	Adjusted PQL	Dilution Factor	PQL	Analytical Method	Analysis Date	By	Prep Method	Prepped Date	By	QC	Notes
ARSENIC	U 0.008	mg/L	0.008	1	0.008	SW846 6010	6/6/06	EAM	SW846 3010	6/2/06	SJC	WF02ICW1	
BARIUM	U 0.0050	mg/L	0.0050	1	0.005	SW846 6010	6/6/06	EAM	SW846 3010	6/2/06	SJC	WF02ICW1	
CADMIUM	U 0.0100	mg/L	0.0100	1	0.01	SW846 6010	6/6/06	EAM	SW846 3010	6/2/06	SJC	WF02ICW1	
CHROMIUM	U 0.015	mg/L	0.015	1	0.015	SW846 6010	6/6/06	EAM	SW846 3010	6/2/06	SJC	WF02ICW1	
LEAD	U 0.005	mg/L	0.005	1	0.005	SW846 6010	6/6/06	EAM	SW846 3010	6/2/06	SJC	WF02ICW1	
MERCURY	U 0.20	ug/L	0.20	1	0.2	SW846 7470	6/5/06	MW	SW846 7470	6/5/06	MW	WF05HGW0	
SELENIUM	U 0.010	mg/L	0.010	1	0.01	SW846 6010	6/6/06	EAM	SW846 3010	6/2/06	SJC	WF02ICW1	
SILVER	U 0.015	mg/L	0.015	1	0.015	SW846 6010	6/6/06	EAM	SW846 3010	6/2/06	SJC	WF02ICW1	

FORM 4

SEMIVOLATILE METHOD BLANK SUMMARY

CLIENT SAMPLE ID

WG28762-BLANK

Lab Name: KATAHDIN ANALYTICAL SERVICES Lab Code: KAS

Project: 218736 SDG No.: WW2676

Lab File ID: X0023 Lab Sample ID: WG28762-1

Instrument ID: GCMS-X Date Extracted: 06/04/06

Matrix: (soil/water) WATER Date Analyzed: 06/06/06

Level:(low/med) LOW Time Analyzed: 1639

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES, MS and MSD:

	CLIENT SAMPLE ID	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED	TIME ANALYZED
01	WG28762-LCS	WG28762-2	X0024	06/06/06	1723
02	WG28762-LCSD	WG28762-3	X0025	06/06/06	1807
03	EB-101	WW2676-1	X0028	06/06/06	2021
04					
05					
06					
07					
08					
09					
10					
11					
12					
13					
14					
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24					
25					
26					
27					
28					
29					
30					

COMMENTS:

FORM IV SV

KATAHDIN ANALYTICAL SERVICES
Report of Analytical Results

Client:
Project: 218736
PO No:
Sample Date:
Received Date:
Extraction Date: 06/04/06
Analysis Date: 06-JUN-2006 16:39
Report Date: 06/07/2006
Matrix: WATER
% Solids: NA

Lab ID: WG28762-1
Client ID: WG28762-Blank
SDG: WW2676
Extracted by: KM
Extraction Method: SW846 3510
Analyst: JCG
Analysis Method: SW846 8270C
Lab Prep Batch: WG28762
Units: ug/L

Compound	Flags	Results	DF	PQL	Adj.PQL
Phenol	U	10	1.0	10	10
Bis(2-Chloroethyl)ether	U	10	1.0	10	10
2-Chlorophenol	U	10	1.0	10	10
1,3-Dichlorobenzene	U	10	1.0	10	10
1,4-Dichlorobenzene	U	10	1.0	10	10
1,2-Dichlorobenzene	U	10	1.0	10	10
2-Methylphenol	U	10	1.0	10	10
2,2'-Oxybis(1-Chloropropane)	U	10	1.0	10	10
N-Nitroso-di-n-propylamine	U	10	1.0	10	10
3&4-Methylphenol	U	10	1.0	10	10
Hexachloroethane	U	10	1.0	10	10
Nitrobenzene	U	10	1.0	10	10
Isophorone	U	10	1.0	10	10
2-Nitrophenol	U	10	1.0	10	10
2,4-Dimethylphenol	U	10	1.0	10	10
Bis(2-Chloroethoxy)methane	U	10	1.0	10	10
2,4-Dichlorophenol	U	10	1.0	10	10
1,2,4-Trichlorobenzene	U	10	1.0	10	10
Naphthalene	U	10	1.0	10	10
4-Chloroaniline	U	10	1.0	10	10
Hexachlorobutadiene	U	10	1.0	10	10
4-Chloro-3-Methylphenol	U	10	1.0	10	10
2-Methylnaphthalene	U	10	1.0	10	10
Hexachlorocyclopentadiene	U	10	1.0	10	10
2,4,6-Trichlorophenol	U	10	1.0	10	10
2,4,5-Trichlorophenol	U	25	1.0	25	25
2-Chloronaphthalene	U	10	1.0	10	10
2-Nitroaniline	U	25	1.0	25	25
Dimethyl Phthalate	U	10	1.0	10	10
2,6-Dinitrotoluene	U	10	1.0	10	10
Acenaphthylene	U	10	1.0	10	10
3-Nitroaniline	U	25	1.0	25	25
Acenaphthene	U	10	1.0	10	10
2,4-Dinitrophenol	U	25	1.0	25	25
Dibenzofuran	U	10	1.0	10	10
4-Nitrophenol	U	25	1.0	25	25
2,4-Dinitrotoluene	U	10	1.0	10	10
Diethylphthalate	U	10	1.0	10	10
Fluorene	U	10	1.0	10	10
4-Chlorophenyl-phenylether	U	10	1.0	10	10
4-Nitroaniline	U	25	1.0	25	25
4,6-Dinitro-2-Methylphenol	U	25	1.0	25	25
N-Nitrosodiphenylamine	U	10	1.0	10	10

KATAHDIN ANALYTICAL SERVICES
Report of Analytical Results

Client:
Project: 218736
PO No:
Sample Date:
Received Date:
Extraction Date: 06/04/06
Analysis Date: 06-JUN-2006 16:39
Report Date: 06/07/2006
Matrix: WATER
% Solids: NA

Lab ID: WG28762-1
Client ID: WG28762-Blank
SDG: WW2676
Extracted by: KM
Extraction Method: SW846 3510
Analyst: JCG
Analysis Method: SW846 8270C
Lab Prep Batch: WG28762
Units: ug/L

Compound	Flags	Results	DF	PQL	Adj.PQL
4-Bromophenyl-phenylether	U	10	1.0	10	10
Hexachlorobenzene	U	10	1.0	10	10
Pentachlorophenol	U	25	1.0	25	25
Phenanthrene	U	10	1.0	10	10
Anthracene	U	10	1.0	10	10
Carbazole	U	10	1.0	10	10
Di-n-butylphthalate	U	10	1.0	10	10
Fluoranthene	U	10	1.0	10	10
Pyrene	U	10	1.0	10	10
Butylbenzylphthalate	U	10	1.0	10	10
Benzo(a)anthracene	U	10	1.0	10	10
3,3'-Dichlorobenzidine	U	10	1.0	10	10
Chrysene	U	10	1.0	10	10
bis(2-Ethylhexyl)phthalate		110	1.0	10	10
Di-n-octylphthalate	U	10	1.0	10	10
Benzo(b)fluoranthene	U	10	1.0	10	10
Benzo(k)fluoranthene	U	10	1.0	10	10
Benzo(a)pyrene	U	10	1.0	10	10
Indeno(1,2,3-cd)pyrene	U	10	1.0	10	10
Dibenzo(a,h)anthracene	U	10	1.0	10	10
Benzo(g,h,i)perylene	U	10	1.0	10	10
2-Fluorophenol		35%			
Phenol-D6		42%			
Nitrobenzene-D5		81%			
2-Fluorobiphenyl		89%			
2,4,6-Tribromophenol		143%			
Terphenyl-D14		114%			

KATAHDIN ANALYTICAL SERVICES
LAB CONTROL SAMPLE

Client:
Project: 218736
PO No:
Sample Date:
Received Date:
Extraction Date: 06/04/06
Analysis Date: 06/06/06
Report Date: 06/07/2006
Matrix: WATER

Lab ID: WG28762-2 & WG28762-3
Client ID: WG28762-LCS & WG28762-LCSD
SDG: WW2676
Extracted by: KM
Extraction Method: SW846 3510
Analyst: JCG
Analysis Method: SW846 8270C
Lab Prep Batch: WG28762
Units: ug/L

COMPOUND	LCS SPIKE	LCSD SPIKE	SAMPLE CONC.	LCS CONC.	LCSD CONC.	LCS %REC.	LCSD %REC.	%RPD	%RPD LIMIT	QC. LIMITS
Phenol	100	100	NA	49	47	49	47	5	30	0-126
Bis(2-Chloroethyl)ether	50	50	NA	40	40	80	80	1	30	35-109
2-Chlorophenol	100	100	NA	91	92	91	92	0.6	30	52-112
1,3-Dichlorobenzene	50	50	NA	35	35	70	70	0.3	30	20-100
1,4-Dichlorobenzene	50	50	NA	36	36	73	73	0.0	30	19-103
1,2-Dichlorobenzene	50	50	NA	38	38	77	77	0.0	30	25- 93
2-Methylphenol	100	100	NA	90	83	90	83	8	30	31-105
2,2'-Oxybis(1-Chloropropane)	50	50	NA	36	34	73	67	8	30	32-108
N-Nitroso-di-n-propylamine	50	50	NA	43	40	85	80	6	30	36-106
3&4-Methylphenol	100	100	NA	78	75	78	75	4	30	24-116
Hexachloroethane	50	50	NA	30	28	61	56	8	30	30- 95
Nitrobenzene	50	50	NA	46	42	91	85	7	30	44-112
Isophorone	50	50	NA	48	45	95	90	5	30	50-110
2-Nitrophenol	100	100	NA	99	98	99	98	0.8	30	49-108
2,4-Dimethylphenol	100	100	NA	96	94	96	94	2	30	48-105
Bis(2-Chloroethoxy)methane	50	50	NA	43	42	85	83	3	30	52-101
2,4-Dichlorophenol	100	100	NA	113	107	* 113	107	5	30	57-110
1,2,4-Trichlorobenzene	50	50	NA	48	45	96	90	6	30	31-114
Naphthalene	50	50	NA	46	43	93	87	6	30	33-109
4-Chloroaniline	50	50	NA	34	36	68	72	5	30	29- 98
Hexachlorobutadiene	50	50	NA	53	46	105	92	13	30	30-107
4-Chloro-3-Methylphenol	100	100	NA	104	103	104	103	1.0	30	46-150
2-Methylnaphthalene	50	50	NA	48	47	97	94	3	30	48-121
Hexachlorocyclopentadiene	50	50	NA	24	20	48	39	20	30	0-106
2,4,6-Trichlorophenol	100	100	NA	115	111	* 115	* 111	4	30	69-105
2,4,5-Trichlorophenol	100	100	NA	95	92	95	92	3	30	58-117
2-Chloronaphthalene	50	50	NA	39	39	78	78	0.5	30	33-119
2-Nitroaniline	50	50	NA	42	41	83	82	0.7	30	45-127
Dimethyl Phthalate	50	50	NA	6.6	6.5	13	13	1	30	4-149
2,6-Dinitrotoluene	50	50	NA	54	51	108	103	4	30	51-120
Acenaphthylene	50	50	NA	46	44	91	88	4	30	49-117
3-Nitroaniline	50	50	NA	38	39	76	78	3	30	20-125
Acenaphthene	50	50	NA	42	40	84	81	4	30	45-118
2,4-Dinitrophenol	100	100	NA	90	104	90	104	15	30	15-140
Dibenzofuran	50	50	NA	46	44	92	88	4	30	53-124
4-Nitrophenol	100	100	NA	64	59	64	59	7	30	0-158
2,4-Dinitrotoluene	50	50	NA	48	47	96	94	1	30	44-133
Diethylphthalate	50	50	NA	31	30	62	61	2	30	20-139
Fluorene	50	50	NA	48	47	96	95	2	30	46-127
4-Chlorophenyl-phenylether	50	50	NA	54	54	109	108	0.6	30	48-136
4-Nitroaniline	50	50	NA	38	39	76	78	3	30	15-145
4,6-Dinitro-2-Methylphenol	100	100	NA	121	123	121	123	2	30	4-150
N-Nitrosodiphenylamine	50	50	NA	45	44	90	87	4	30	30-150
4-Bromophenyl-phenylether	50	50	NA	68	70	* 136	* 139	2	30	68-129
Hexachlorobenzene	50	50	NA	82	81	* 165	* 163	1	30	67-141

KATAHDIN ANALYTICAL SERVICES
LAB CONTROL SAMPLE

Client:
Project: 218736
PO No:
Sample Date:
Received Date:
Extraction Date: 06/04/06
Analysis Date: 06/06/06
Report Date: 06/07/2006
Matrix: WATER

Lab ID: WG28762-2 & WG28762-3
Client ID: WG28762-LCS & WG28762-LCSD
SDG: WW2676
Extracted by: KM
Extraction Method: SW846 3510
Analyst: JCG
Analysis Method: SW846 8270C
Lab Prep Batch: WG28762
Units: ug/L

COMPOUND	LCS SPIKE	LCSD SPIKE	SAMPLE CONC.	LCS CONC.	LCSD CONC.	LCS %REC.	LCSD %REC.	%RPD	%RPD LIMIT	QC. LIMITS
Pentachlorophenol	100	100	NA	158	146	158	146	8	30	10-174
Phenanthrene	50	50	NA	56	58	113	115	2	30	51-139
Anthracene	50	50	NA	57	58	114	115	0.9	30	53-133
Carbazole	50	50	NA	47	47	94	94	0.0	30	35-157
Di-n-butylphthalate	50	50	NA	48	46	96	92	4	30	14-165
Fluoranthene	50	50	NA	54	56	109	112	3	30	32-172
Pyrene	50	50	NA	68	60	* 137	* 121	12	30	49-119
Butylbenzylphthalate	50	50	NA	34	32	68	64	6	30	30-126
Benzo(a)anthracene	50	50	NA	56	56	* 113	* 112	0.4	30	53-111
3,3'-Dichlorobenzidine	50	50	NA	45	43	90	86	4	30	38-101
Chrysene	50	50	NA	60	57	120	114	6	30	59-120
bis(2-Ethylhexyl)phthalate	50	50	NA	54	51	107	101	6	30	39-125
Di-n-octylphthalate	50	50	NA	49	50	98	99	0.6	30	31-136
Benzo(b)fluoranthene	50	50	NA	52	49	104	99	5	30	52-120
Benzo(k)fluoranthene	50	50	NA	52	50	105	99	6	30	56-126
Benzo(a)pyrene	50	50	NA	50	49	99	98	1	30	51-120
Indeno(1,2,3-cd)pyrene	50	50	NA	48	45	96	90	7	30	32-143
Dibenzo(a,h)anthracene	50	50	NA	48	47	97	95	2	30	37-134
Benzo(g,h,i)perylene	50	50	NA	47	44	94	88	7	30	32-143

FORM III SV-1

X0024.D & X0025.D

FORM 4
PESTICIDE METHOD BLANK SUMMARY

CLIENT SAMPLE ID

WG28773-BLANK

Lab Name: KATAHDIN ANALYTICAL SERVICES Lab Code: KAS

Project: 218736 SDG No.: WW2676

Lab Sample ID: WG28773-1 Lab File ID: 6WF1178

Matrix (soil/water) WATER Extraction:(SepF/Cont/Sonc) SW846 3510

Sulfur Cleanup: (Y/N) N Date Extracted: 06/05/06

Date Analyzed (1): 06/07/06 Date Analyzed (2): 06/07/06

Time Analyzed (1): 0144 Time Analyzed (2): 0144

Instrument ID (1): GC06 Instrument ID (2): GC06

GC Column (1): RTX-5 ID: 0.53(mm) GC Column (2): RTX-35 ID: 0.53(mm)

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES, MS and MSD:

	CLIENT SAMPLE ID	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED 1	DATE ANALYZED 2
01	EB-101	WW2676-1	6WF1179	06/07/06	06/07/06
02	WG28773-LCS	WG28773-2	6WF1199	06/07/06	06/07/06
03					
04					
05					
06					
07					
08					
09					
10					
11					
12					
13					
14					
15					
16					
17					
18					
19					
20					
21					
22					
23					
24					

COMMENTS: _____

FORM IV PEST

KATAHDIN ANALYTICAL SERVICES
Report of Analytical Results

Client:
Project: 218736
PO No:
Sample Date:
Received Date:
Extraction Date: 06/05/06
Analysis Date: 07-JUN-2006 01:44
Report Date: 06/07/2006
Matrix: WATER
% Solids: NA

Lab ID: WG28773-1
Client ID: WG28773-Blank
SDG: WW2676
Extracted by: KM
Extraction Method: SW846 3510
Analyst: JLP
Analysis Method: SW846 8082
Lab Prep Batch: WG28773
Units: ug/L

Compound	Flags	Results	DF	PQL	Adj.PQL
Aroclor-1016	U	0.50	1.0	0.50	0.50
Aroclor-1221	U	0.50	1.0	0.50	0.50
Aroclor-1232	U	0.50	1.0	0.50	0.50
Aroclor-1242	U	0.50	1.0	0.50	0.50
Aroclor-1248	U	0.50	1.0	0.50	0.50
Aroclor-1254	U	0.50	1.0	0.50	0.50
Aroclor-1260	U	0.50	1.0	0.50	0.50
Tetrachloro-m-xylene		70%			
Decachlorobiphenyl		57%			

KATAHDIN ANALYTICAL SERVICES
LAB CONTROL SAMPLE

Client:
Project: 218736
PO No:
Sample Date:
Received Date:
Extraction Date: 06/05/06
Analysis Date: 06/07/06
Report Date: 06/07/2006
Matrix: WATER

Lab ID: WG28773-2
Client ID: WG28773-LCS
SDG: WW2676
Extracted by: KM
Extraction Method: SW846 3510
Analyst: JLP
Analysis Method: SW846 8082
Lab Prep Batch: WG28773
Units: ug/L

COMPOUND	LCS SPIKE	SAMPLE CONC.	LCS CONC.	%REC.	QC. LIMITS
Aroclor-1016	5.0	NA	5.0	100	72-119
Aroclor-1260	5.0	NA	4.3	87	60-119

Katahdin
ANALYTICAL SERVICES

PREPARATION BLANK REPORT

Sample ID: PBWWF02ICW1 **Batch ID** WF02ICW1

Element Name	Result	Units	Flag	PQL	File
ALUMINUM	0.019	mg/L	J	0.300	BWF02A
ANTIMONY	0.004	mg/L	U	0.008	AWF06B
ARSENIC	0.003	mg/L	U	0.008	AWF06B
BARIUM	0.001	mg/L	U	0.005	BWF02A
BERYLLIUM	0.0003	mg/L	U	0.0050	BWF02A
BORON	0.007	mg/L	U	0.100	BWF02A
CADMIUM	0.004	mg/L	U	0.010	BWF02A
CALCIUM	0.02	mg/L	U	0.05	BWF06A
CHROMIUM	0.001	mg/L	U	0.015	AWF06B
COBALT	0.003	mg/L	U	0.030	BWF02A
COPPER	0.003	mg/L	U	0.025	BWF02A
IRON	0.007	mg/L	J	0.100	BWF02A
LEAD	0.002	mg/L	U	0.005	AWF06B
MAGNESIUM	0.03	mg/L	U	0.05	BWF02A
MANGANESE	0.0007	mg/L	U	0.0050	BWF02A
MOLYBDENUM	0.002	mg/L	U	0.010	AWF06B
NICKEL	0.01	mg/L	U	0.04	BWF02A
POTASSIUM	0.4	mg/L	U	1.0	BWF02A
SELENIUM	0.004	mg/L	U	0.010	AWF06B
SILICON	0.05	mg/L	J	0.20	BWF02A
SILVER	0.004	mg/L	U	0.015	BWF02A
SODIUM	0.1	mg/L	J	1.00	BWF06A
STRONTIUM	0.00005	mg/L	U	0.100	AWF06B
TIN	0.003	mg/L	U	0.100	AWF06B
VANADIUM	0.005	mg/L	U	0.025	BWF02A
ZINC	0.002	mg/L	U	0.025	BWF02A

U The analyte was not detected in the sample at a level greater than the instrument detection limit.

J The analyte was detected in the sample at a concentration greater than the instrument detection limit, but less than the laboratory's Practical Quantitation Level.

H The analyte was detected in the sample at a concentration greater than the laboratory's acceptance limit.

Katahdin
ANALYTICAL SERVICES

LABORATORY CONTROL SAMPLE REPORT

Sample ID: LCSWWF02ICW1 **Batch ID** WF02ICW1

Element Name	True Value	Result	Units	Recovery(%)	Flag	Limits (%)		File
ALUMINUM	2.00	2.08	mg/L	104.0%		80.	120.	BWF02A
ANTIMONY	0.500	0.468	mg/L	93.6%		80.	120.	AWF06B
ARSENIC	0.500	0.459	mg/L	91.8%		80.	120.	AWF06B
BARIUM	2.00	2.12	mg/L	106.0%		80.	120.	BWF02A
BERYLLIUM	0.0500	0.0508	mg/L	101.6%		80.	120.	BWF02A
BORON	0.500	0.516	mg/L	103.2%		80.	120.	BWF02A
CADMIUM	0.250	0.258	mg/L	103.2%		80.	120.	BWF02A
CALCIUM	2.50	2.76	mg/L	110.4%		80.	120.	BWF06A
CHROMIUM	0.200	0.216	mg/L	108.0%		80.	120.	AWF06B
COBALT	0.500	0.525	mg/L	105.0%		80.	120.	BWF02A
COPPER	0.250	0.256	mg/L	102.4%		80.	120.	BWF02A
IRON	1.00	1.06	mg/L	106.0%		80.	120.	BWF02A
LEAD	0.500	0.475	mg/L	95.0%		80.	120.	AWF06B
MAGNESIUM	5.00	5.28	mg/L	105.6%		80.	120.	BWF02A
MANGANESE	0.500	0.517	mg/L	103.4%		80.	120.	BWF02A
MOLYBDENUM	0.300	0.293	mg/L	97.7%		80.	120.	AWF06B
NICKEL	0.50	0.52	mg/L	104.0%		80.	120.	BWF02A
POTASSIUM	10.0	10.8	mg/L	108.0%		80.	120.	BWF02A
SELENIUM	0.500	0.456	mg/L	91.2%		80.	120.	AWF06B
SILICON	0.98	1.09	mg/L	111.2%		80.	120.	BWF02A
SILVER	0.050	0.046	mg/L	92.0%		80.	120.	BWF02A
SODIUM	7.50	8.23	mg/L	109.7%		80.	120.	BWF02A
STRONTIUM	0.750	0.781	mg/L	104.1%		80.	120.	AWF06B
TIN	0.750	0.705	mg/L	94.0%		80.	120.	AWF06B
VANADIUM	0.500	0.523	mg/L	104.6%		80.	120.	BWF02A
ZINC	0.500	0.524	mg/L	104.8%		80.	120.	BWF02A

H Laboratory control sample recovery is greater than the laboratory's acceptance limit.

L Laboratory control sample recovery is less than the laboratory's acceptance limit.

Katahdin
ANALYTICAL SERVICES

PREPARATION BLANK REPORT

Sample ID: PBWWF05HGW0 **Batch ID** WF05HGW0

Element Name	Result	Units	Flag	PQL	File
MERCURY	0.02	ug/L	U	0.20	HWF05A

U The analyte was not detected in the sample at a level greater than the instrument detection limit.

J The analyte was detected in the sample at a concentration greater than the instrument detection limit, but less than the laboratory's Practical Quantitation Level.

H The analyte was detected in the sample at a concentration greater than the laboratory's acceptance limit.

Katahdin
ANALYTICAL SERVICES

LABORATORY CONTROL SAMPLE REPORT

Sample ID: LCSWWF05HGW0 **Batch ID** WF05HGW0

Element Name	True Value	Result	Units	Recovery(%)	Flag	Limits (%)		File
MERCURY	5.00	4.83	ug/L	96.6%		80.	120.	HWF05A

H Laboratory control sample recovery is greater than the laboratory's acceptance limit.

L Laboratory control sample recovery is less than the laboratory's acceptance limit.

KATAHDIN ANALYTICAL SERVICES, INC.

SAMPLE RECEIPT CONDITION REPORT

Tel. (207) 874-2400
Fax (207) 775-4029

LAB (WORK ORDER) #: _____

PAGE: 1 OF 1

COOLER: 1 OF 1

CLIENT: W&C

PROJECT: _____

COC#: _____
SDG#: _____
DATE / TIME RECEIVED: 6-1-06 1310
DELIVERED BY: Client
RECEIVED BY: D. Warneke
LIMS ENTRY BY: lwm
LIMS REVIEW BY / PM: AJC

	YES	NO	EXCEPTIONS	COMMENTS	RESOLUTION
1. CUSTODY SEALS PRESENT / INTACT?	☒	☐	☐		
2. CHAIN OF CUSTODY PRESENT IN THIS COOLER?	☒	☐	☐		
3. CHAIN OF CUSTODY SIGNED BY CLIENT?	☒	☐	☐		
4. CHAIN OF CUSTODY MATCHES SAMPLES?	☒	☐	☐		
5. TEMPERATURE BLANKS PRESENT?	☒	☐	☐	TEMP BLANK TEMP (°C)= 14.0	
6. SAMPLES RECEIVED AT 4°C +/- 2? ICE / ICE PACKS PRESENT (Y) or N?	☐	☒	☐	COOLER TEMP (°C)= 20.0 NA (RECORD COOLER TEMP ONLY IF TEMP BLANK IS NOT PRESENT)	
7. VOLATILES FREE OF HEADSPACE?	☐	☐	☒		
8. TRIP BLANK PRESENT IN THIS COOLER	☐	☐	☒		
9. PROPER SAMPLE CONTAINERS AND VOLUME?	☒	☐	☐		
10. SAMPLES WITHIN HOLD TIME UPON RECEIPT?	☒	☐	☐		
11. SAMPLES PROPERLY PRESERVED[1]?	☒	☐	☐		
12. CORRECTIVE ACTION REPORT FILED?	☐	☒	N/A		

13. ANALYTICAL PROGRAMS (CIRCLE ONE) COMMERCIAL CLP HAZWRAP NFESC ACOE AFCEE OTHER (STATE OF ORIGIN): _____

LOG - IN NOTES[1]:

[1] Use this space (and additional sheets if necessary) to document samples that are received broken or compromised, C-O-C discrepancies, radiation checks, residual chlorine check, results of pH check if required. If samples required pH adjustment, record volume and type of preservative added.

Katahdin Analytical Services WW2676 page 0000020 of 0000022

Katahdin ANALYTICAL SERVICES
340 County Road No. 5
P.O. Box 720
Westbrook, ME 04092
Tel: (207) 874-2400
Fax: (207) 775-4029

CHAIN of CUSTODY

PLEASE BEAR DOWN AND PRINT LEGIBLY IN PEN

Client	Contact	Phone #	Fax #
WOODARD & CURRAN	KARL KASPER	(207) 774-2112	( )

Address	City	State	Zip Code
41 HUTCHINS DR	PORTLAND	ME	04103

Purchase Order #	Proj. Name / No.	Katahdin Quote #
	218736	

Bill (if different than above) Address

Sampler (Print / Sign)

Copies To:

LAB USE ONLY WORK ORDER #: WW2676

KATAHDIN PROJECT NUMBER

REMARKS:

SHIPPING INFO: ☐ FED EX ☐ UPS ☐ CLIENT

AIRBILL NO:

TEMP°C ☐ TEMP BLANK ☐ INTACT ☐ NOT INTACT

ANALYSIS AND CONTAINER TYPE PRESERVATIVES

*	Sample Description	Date / Time coll'd	Matrix	No. of Cntrs.	RCRA8 metals Filt. ☐Y ☒N	SVOC's Filt. ☐Y ☒N	~~Water~~ PCBs Filt. ☐Y ☒N	Filt. ☐Y ☐N	Filt. ☐Y ☐N	Filt. ☐Y ☐N	Filt. ☐Y ☐N	Filt. ☐Y ☐N	Filt. ☐Y ☐N	Filt. ☐Y ☐N
	EB-101	6/01/06 1230	GW	3	1	1	1							
		/												
		/												
		/												
		/												
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		/												
		/												

COMMENTS

RESULTS BY NEXT WEDNESDAY 7 JUNE – ANALYTE IF OUT OF TEMP RANGE

Relinquished By: (Signature)	Date / Time	Received By: (Signature)	Relinquished By: (Signature)	Date / Time	Received By: (Signature)
	6/01/06 1250				
Relinquished By: (Signature)	Date / Time	Received By: (Signature)	Relinquished By: (Signature)	Date / Time	Received By: (Signature)

THE TERMS AND CONDITIONS ON THE REVERSE SIDE HEREOF SHALL GOVERN SERVICES, EXCEPT WHEN A SIGNED CONTRACTUAL AGREEMENT EXISTS.

ORIGINAL

Katahdin
ANALYTICAL SERVICES

Katahdin Analytical Services

Login Chain of Custody Report (lno1)

Jun. 01, 2006
05:21 PM

Login Number: WW2676

Account: WOODAR001 Web
Woodard & Curran

Project:

Primary Report Address:
Erika Lloyd
Woodard & Curran
41 Hutchins Drive

Portland,ME 04102

Primary Invoice Address:
Accounts Payable
Woodard & Curran
41 Hutchins Drive

Portland,ME 04102

Report CC Addresses:
Invoice CC Addresses:

Login Information

ANALYSIS INSTRUCTIONS	:
CHECK NO.	:
CLIENT PO#	:
COOLER TEMPERATURE	: 14.0
DELIVERY SERVICES	: Client
EDD FORMAT	: KAS027-XLS & KAS064-XLS
MAIL DATE	:
PM	: AJC
PROJECT NAME	: 218736
QC LEVEL	: II
REGULATORY LIST	:
REPORT INSTRUCTIONS	:
SDG ID	:
SDG STATUS	:

Laboratory Sample ID	Client Sample Number	Collect Date/Time	Receive Date	PR	Verbal Date	Due Date	Comments
WW2676-1	EB-101	01-JUN-06 12:30	01-JUN-06		**07-JUN-06**	07-JUN-06	

Matrix		*Product*	*Hold Date (shortest)*	*Bottle Type*	*Bottle Count*
Aqueous	P	RCRA-METALS			1
SW3010-PREP		SW6010-ARSENIC	SW6010-BARIUM	SW6010-CADMIUM	
SW6010-CHROMIUM		SW6010-LEAD	SW6010-SELENIUM	SW6010-SILVER	
SW7470-MERCURY					
Aqueous	S	SW8082	08-JUN-06	1L N-Amber Glass	1
Aqueous	S	SW8270BNA	08-JUN-06	1L N-Amber Glass	1

Total Samples: *1* ***Total Analyses:*** *3*

APPENDIX E: SUBJECT PROPERTY PHOTOLOG

Photo 1. SB-3.

Photo 2. SB-3.

Photo 3. SB-10.

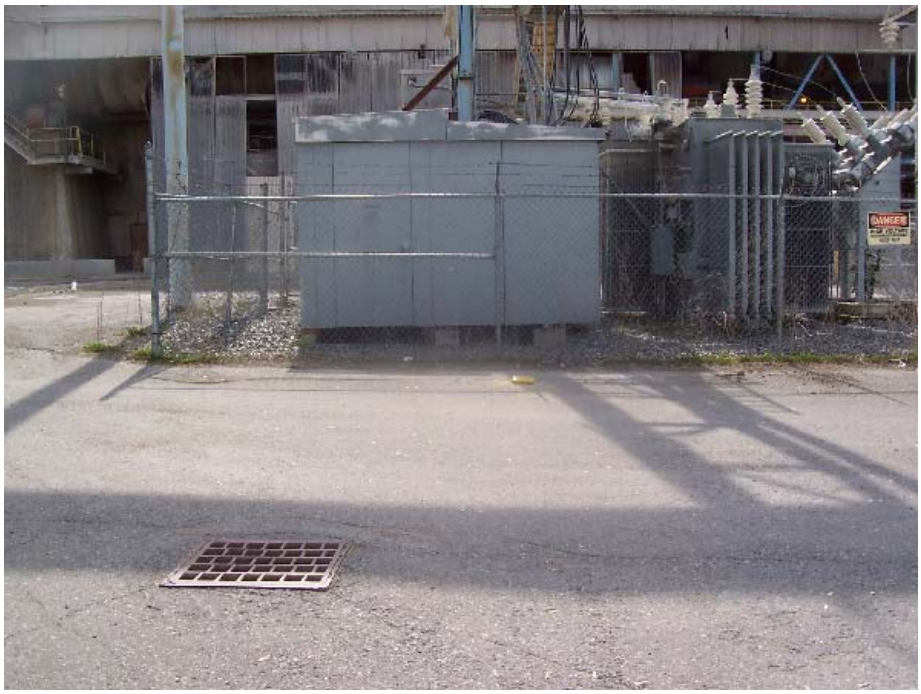

Photo 4. SB-10.

Photo 5. SB-13.

Photo 6. SB-15.

Photo 7. SB-15.

Photo 8. SB-18.

Photo 9. SB-19.

Photo 10. SB-20.

Photo 11. SB-30.

Photo 12. SB-30.

Photo 13. SB-37.

Photo 14. SB-37.

Photo 15. SB-38.

Photo 16. SB-38.

Photo 17. Former South Bark Yard.

Photo 18. Former South Bark Yard.

Photo 19. Former South Bark Yard.

Photo 20. Former South Bark Yard.

Photo 21. Former South Bark Yard.

Photo 22. TP-101.

Photo 23. TP-108.

Photo 24. TP-121.