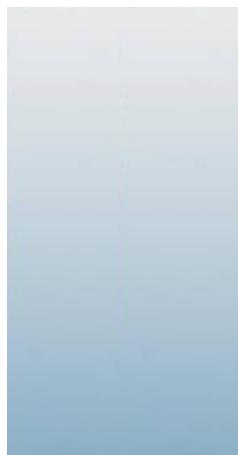




11/15/12



## Technical Report for

**XTO Energy**

**PCU 296-6A**

**1211-02**

**Accutest Job Number: D40798**

**Sampling Date: 11/08/12**

### Report to:

**KRW Consulting, Inc.  
8000 West 14th Avenue  
Lakewood, CO 80214  
dknudson@krwconsulting.com; jhess@krwconsulting.com;  
crachak@krwconsulting.com; rrasic@krwconsulting.com;  
ATTN: Dwayne Knudson**

**Total number of pages in report: 143**



Test results contained within this data package meet the requirements of the National Environmental Laboratory Accreditation Conference and/or state specific certification programs as applicable.

A handwritten signature in black ink.

**Brad Madadian  
Laboratory Director**

**Client Service contact: Renea Jackson 303-425-6021**

Certifications: CO, ID, NE, NM, ND (R-027) (PW), UT (NELAP CO00049), TX (T104704511-12-1)

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Test results relate only to samples analyzed.

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## Sample Summary

**XTO Energy****Job No:** D40798**PCU 296-6A****Project No:** 1211-02

Sample Number	Collected Date	Time By	Matrix Received	Code Type	Client Sample ID
D40798-1	11/08/12	11:50 DS	11/10/12	SO Soil	CUT 2 SUBLINER (COMP)
D40798-1A	11/08/12	11:50 DS	11/10/12	SO Soil	CUT 2 SUBLINER (COMP)

---

Soil samples reported on a dry weight basis unless otherwise indicated on result page.



## CASE NARRATIVE / CONFORMANCE SUMMARY

**Client:** XTO Energy

**Job No** D40798

**Site:** PCU 296-6A

**Report Date** 11/15/2012 1:48:00 PM

On 11/10/2012, 1 sample(s), 0 Trip Blank(s), and 0 Field Blank(s) were received at Accutest Mountain States (AMS) at a temperature of 3.5 °C. The samples were intact and properly preserved, unless noted below. An AMS Job Number of D40798 was assigned to the project. The lab sample ID, client sample ID, and date of sample collection are detailed in the report's Results Summary.

Specified quality control criteria were achieved for this job except as noted below. For more information, please refer to the analytical results and QC summary pages.

### Volatiles by GCMS By Method SW846 8260B

<b>Matrix</b> SO	<b>Batch ID:</b> V5V1501
------------------	--------------------------

- All samples were analyzed within the recommended method holding time.
- Sample(s) D40713-1MS, D40713-1MSD were used as the QC samples indicated.
- All method blanks for this batch meet method specific criteria.

### Extractables by GCMS By Method SW846 8270C BY SIM

<b>Matrix</b> SO	<b>Batch ID:</b> OP6947
------------------	-------------------------

- All samples were extracted within the recommended method holding time.
- All samples were analyzed within the recommended method holding time.
- All method blanks for this batch meet method specific criteria.
- Sample(s) D40799-1MS, D40799-1MSD were used as the QC samples indicated.

### Volatiles by GC By Method SW846 8015B

<b>Matrix</b> SO	<b>Batch ID:</b> GGB1006
------------------	--------------------------

- All samples were analyzed within the recommended method holding time.
- All method blanks for this batch meet method specific criteria.
- Sample(s) D40795-1MS, D40795-1MSD were used as the QC samples indicated.

### Extractables by GC By Method SW846-8015B

<b>Matrix</b> SO	<b>Batch ID:</b> OP6948
------------------	-------------------------

- All samples were extracted within the recommended method holding time.
- All samples were analyzed within the recommended method holding time.
- Sample(s) D40799-1MS, D40799-1MSD were used as the QC samples indicated.
- All method blanks for this batch meet method specific criteria.

## Metals By Method SW846 6010C

**Matrix** AQ

**Batch ID:** MP8877

- All samples were digested within the recommended method holding time.
- All samples were analyzed within the recommended method holding time.
- All method blanks for this batch meet method specific criteria.
- Sample(s) D40797-1AMS, D40797-1AMSD, D40797-1ASDL were used as the QC samples for the metals analysis.

**Matrix** SO

**Batch ID:** MP8869

- All samples were digested within the recommended method holding time.
- All samples were analyzed within the recommended method holding time.
- All method blanks for this batch meet method specific criteria.
- Sample(s) D40778-1MS, D40778-1MSD, D40778-1SDL were used as the QC samples for the metals analysis.
- The serial dilution RPD(s) for Selenium, Silver, Barium, Chromium, Nickel, Zinc are outside control limits for sample MP8869-SD1. Percent difference acceptable due to low initial sample concentration (< 50 times IDL).
- MP8869-SD1 for Chromium: Serial dilution indicates possible matrix interference.
- MP8869-SD1 for Nickel: Serial dilution indicates possible matrix interference.
- MP8869-SD1 for Zinc: Serial dilution indicates possible matrix interference.
- MP8869-SD1 for Barium: Serial dilution indicates possible matrix interference.

## Metals By Method SW846 6020A

**Matrix** SO

**Batch ID:** MP8870

- All samples were digested within the recommended method holding time.
- All samples were analyzed within the recommended method holding time.
- All method blanks for this batch meet method specific criteria.
- Sample(s) D40778-1MS, D40778-1MSD, D40778-1SDL were used as the QC samples for the metals analysis.

## Metals By Method SW846 7471B

**Matrix** SO

**Batch ID:** MP8871

- All samples were digested within the recommended method holding time.
- All samples were analyzed within the recommended method holding time.
- All method blanks for this batch meet method specific criteria.
- Sample(s) D40797-1MS, D40797-1MSD were used as the QC samples for the metals analysis.

## Wet Chemistry By Method ASTM D1498-76M

**Matrix** SO

**Batch ID:** GN17680

- Sample(s) D40797-1DUP were used as the QC samples for the Redox Potential Vs H<sub>2</sub> analysis.

## Wet Chemistry By Method SM19 2540B M

**Matrix** SO

**Batch ID:** GN17634

- The data for SM19 2540B M meets quality control requirements.

## Wet Chemistry By Method SW846 3060/7196A M

**Matrix** SO

**Batch ID:** R15135

- The data for SW846 3060/7196A M meets quality control requirements.
- D40798-1 for Chromium, Trivalent: Calculated as: (Chromium) - (Chromium, Hexavalent)

## Wet Chemistry By Method SW846 3060A/7196A

**Matrix** SO

**Batch ID:** GP8655

- All samples were prepared within the recommended method holding time.
- All samples were analyzed within the recommended method holding time.
- All method blanks for this batch meet method specific criteria.
- Sample(s) D40715-1DUP, D40715-1MS, D40715-1MSD were used as the QC samples for the Chromium, Hexavalent analysis.

## Wet Chemistry By Method SW846 9045D

**Matrix** SO

**Batch ID:** GN17678

- The following samples were run outside of holding time for method SW846 9045D: D40798-1

## Wet Chemistry By Method USDA HANDBOOK 60

**Matrix** SO

**Batch ID:** MP8877

- D40798-1A for Sodium Adsorption Ratio: Calculated as:  $(\text{Na meq/L}) / \sqrt{[(\text{Ca meq/L}) + (\text{Mg meq/L})/2]}$

AMS certifies that data reported for samples received, listed on the associated custody chain or analytical task order, were produced to specifications meeting AMS's Quality System precision, accuracy and completeness objectives except as noted.

Estimated non-standard method measurement uncertainty data is available on request, based on quality control bias and implicit for standard methods. Acceptable uncertainty requires tested parameter quality control data to meet method criteria.

AMS is not responsible for data quality assumptions if partial reports are used and recommends that this report be used in its entirety. This report is authorized by AMS indicated via signature on the report cover.

## Summary of Hits

Page 1 of 1

Job Number: D40798  
Account: XTO Energy  
Project: PCU 296-6A  
Collected: 11/08/12

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Lab Sample ID	Client Sample ID	Result/ Analyte	Qual	RL	MDL	Units	Method
---------------	------------------	--------------------	------	----	-----	-------	--------

### D40798-1 CUT 2 SUBLINER (COMP)

Arsenic	6.4	0.11	mg/kg	SW846 6020A
Barium	438	1.1	mg/kg	SW846 6010C
Chromium	41.8	1.1	mg/kg	SW846 6010C
Copper	7.3	1.1	mg/kg	SW846 6010C
Lead	7.7	5.5	mg/kg	SW846 6010C
Nickel	13.9	3.3	mg/kg	SW846 6010C
Zinc	40.0	3.3	mg/kg	SW846 6010C
Specific Conductivity	377	1.0	umhos/cm	SM2510B-1997 MOD
Chromium, Trivalent <sup>a</sup>	41.8	2.1	mg/kg	SW846 3060/7196A M
Redox Potential Vs H2	141		mv	ASTM D1498-76M
pH	9.75		su	SW846 9045D

### D40798-1A CUT 2 SUBLINER (COMP)

Calcium	15.3	2.0	mg/l	SW846 6010C
Magnesium	4.56	1.0	mg/l	SW846 6010C
Sodium	59.2	2.0	mg/l	SW846 6010C
Sodium Adsorption Ratio <sup>b</sup>	3.41		ratio	USDA HANDBOOK 60

(a) Calculated as: (Chromium) - (Chromium, Hexavalent)

(b) Calculated as: (Na meq/L) / sqrt [(Ca meq/L)+(Mg meq/L)/2]



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## Sample Results

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### Report of Analysis

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## Report of Analysis

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Client Sample ID: CUT 2 SUBLINER (COMP)

Lab Sample ID: D40798-1

Date Sampled: 11/08/12

Matrix: SO - Soil

Date Received: 11/10/12

Method: SW846 8260B

Percent Solids: 89.6

Project: PCU 296-6A

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	5V24609.D	1	11/14/12	BD	n/a	n/a	V5V1501
Run #2							

	Initial Weight	Final Volume	Methanol Aliquot
Run #1	5.02 g	5.0 ml	100 ul
Run #2			

## Purgeable Aromatics

CAS No.	Compound	Result	RL	MDL	Units	Q
71-43-2	Benzene	ND	0.061	0.031	mg/kg	
108-88-3	Toluene	ND	0.12	0.061	mg/kg	
100-41-4	Ethylbenzene	ND	0.12	0.023	mg/kg	
1330-20-7	Xylene (total)	ND	0.25	0.12	mg/kg	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
2037-26-5	Toluene-D8	103%		64-130%
460-00-4	4-Bromofluorobenzene	101%		62-131%
17060-07-0	1,2-Dichloroethane-D4	90%		70-130%

ND = Not detected MDL - Method Detection Limit

J = Indicates an estimated value

RL = Reporting Limit

B = Indicates analyte found in associated method blank

E = Indicates value exceeds calibration range

N = Indicates presumptive evidence of a compound

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## Report of Analysis

Page 1 of 1

Client Sample ID: CUT 2 SUBLINER (COMP)

Lab Sample ID: D40798-1

Date Sampled: 11/08/12

Matrix: SO - Soil

Date Received: 11/10/12

Method: SW846 8270C BY SIM SW846 3546

Percent Solids: 89.6

Project: PCU 296-6A

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	3G12033.D	1	11/12/12	DC	11/12/12	OP6947	E3G568
Run #2							

	Initial Weight	Final Volume
Run #1	30.1 g	1.0 ml
Run #2		

## COGCC Table 910-1 PAH List

CAS No.	Compound	Result	RL	MDL	Units	Q
83-32-9	Acenaphthene	ND	0.0093	0.0048	mg/kg	
120-12-7	Anthracene	ND	0.0093	0.0048	mg/kg	
56-55-3	Benzo(a)anthracene	ND	0.0093	0.0048	mg/kg	
205-99-2	Benzo(b)fluoranthene	ND	0.0093	0.0048	mg/kg	
207-08-9	Benzo(k)fluoranthene	ND	0.0093	0.0048	mg/kg	
50-32-8	Benzo(a)pyrene	ND	0.0093	0.0048	mg/kg	
218-01-9	Chrysene	ND	0.0093	0.0048	mg/kg	
53-70-3	Dibenz(a,h)anthracene	ND	0.0093	0.0048	mg/kg	
206-44-0	Fluoranthene	ND	0.0093	0.0048	mg/kg	
86-73-7	Fluorene	ND	0.0093	0.0048	mg/kg	
193-39-5	Indeno(1,2,3-cd)pyrene	ND	0.0093	0.0048	mg/kg	
91-20-3	Naphthalene	ND	0.013	0.011	mg/kg	
129-00-0	Pyrene	ND	0.0093	0.0048	mg/kg	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
4165-60-0	Nitrobenzene-d5	74%		10-159%
321-60-8	2-Fluorobiphenyl	65%		19-131%
1718-51-0	Terphenyl-d14	79%		18-150%

ND = Not detected MDL - Method Detection Limit

J = Indicates an estimated value

RL = Reporting Limit

B = Indicates analyte found in associated method blank

E = Indicates value exceeds calibration range

N = Indicates presumptive evidence of a compound

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## Report of Analysis

Page 1 of 1

Client Sample ID: CUT 2 SUBLINER (COMP)

Lab Sample ID: D40798-1

Date Sampled: 11/08/12

Matrix: SO - Soil

Date Received: 11/10/12

Method: SW846 8015B

Percent Solids: 89.6

Project: PCU 296-6A

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	GB18417.D	1	11/12/12	SK	n/a	n/a	GGB1006
Run #2							

	Initial Weight	Final Volume	Methanol Aliquot
Run #1	5.0 g	5.0 ml	100 ul
Run #2			

CAS No.	Compound	Result	RL	MDL	Units	Q
	TPH-GRO (C6-C10)	ND	12	6.1	mg/kg	
CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits		
120-82-1	1,2,4-Trichlorobenzene	99%		60-140%		

ND = Not detected MDL - Method Detection Limit

J = Indicates an estimated value

RL = Reporting Limit

B = Indicates analyte found in associated method blank

E = Indicates value exceeds calibration range

N = Indicates presumptive evidence of a compound

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## Report of Analysis

Page 1 of 1

Client Sample ID: CUT 2 SUBLINER (COMP)

Lab Sample ID: D40798-1

Date Sampled: 11/08/12

Matrix: SO - Soil

Date Received: 11/10/12

Method: SW846-8015B SW846 3546

Percent Solids: 89.6

Project: PCU 296-6A

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	FD19426.D	1	11/12/12	AV	11/12/12	OP6948	GFD976
Run #2							

	Initial Weight	Final Volume
Run #1	30.1 g	2.0 ml
Run #2		

CAS No.	Compound	Result	RL	MDL	Units	Q
	TPH-DRO (C10-C28)	ND	15	9.7	mg/kg	
CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits		
84-15-1	o-Terphenyl	91%			35-130%	

ND = Not detected      MDL - Method Detection Limit  
 RL = Reporting Limit  
 E = Indicates value exceeds calibration range

J = Indicates an estimated value  
 B = Indicates analyte found in associated method blank  
 N = Indicates presumptive evidence of a compound

## Report of Analysis

Page 1 of 1

Client Sample ID: CUT 2 SUBLINER (COMP)

Lab Sample ID: D40798-1

Matrix: SO - Soil

Date Sampled: 11/08/12

Date Received: 11/10/12

Percent Solids: 89.6

Project: PCU 296-6A

## Metals Analysis

Analyte	Result	RL	Units	DF	Prep	Analyzed By	Method	Prep Method
Arsenic	6.4	0.11	mg/kg	5	11/12/12	11/15/12	JB	SW846 6020A <sup>3</sup>
Barium	438	1.1	mg/kg	1	11/12/12	11/13/12	JM	SW846 6010C <sup>2</sup>
Cadmium	< 1.1	1.1	mg/kg	1	11/12/12	11/13/12	JM	SW846 6010C <sup>2</sup>
Chromium	41.8	1.1	mg/kg	1	11/12/12	11/13/12	JM	SW846 6010C <sup>2</sup>
Copper	7.3	1.1	mg/kg	1	11/12/12	11/13/12	JM	SW846 6010C <sup>2</sup>
Lead	7.7	5.5	mg/kg	1	11/12/12	11/13/12	JM	SW846 6010C <sup>2</sup>
Mercury	< 0.091	0.091	mg/kg	1	11/13/12	11/13/12	JM	SW846 7471B <sup>1</sup>
Nickel	13.9	3.3	mg/kg	1	11/12/12	11/13/12	JM	SW846 6010C <sup>2</sup>
Selenium	< 5.5	5.5	mg/kg	1	11/12/12	11/13/12	JM	SW846 6010C <sup>2</sup>
Silver	< 3.3	3.3	mg/kg	1	11/12/12	11/13/12	JM	SW846 6010C <sup>2</sup>
Zinc	40.0	3.3	mg/kg	1	11/12/12	11/13/12	JM	SW846 6010C <sup>2</sup>

- (1) Instrument QC Batch: MA2991
- (2) Instrument QC Batch: MA2995
- (3) Instrument QC Batch: MA3000
- (4) Prep QC Batch: MP8869
- (5) Prep QC Batch: MP8870
- (6) Prep QC Batch: MP8871

RL = Reporting Limit

## Report of Analysis

Page 1 of 1

Client Sample ID:	CUT 2 SUBLINER (COMP)	Date Sampled:	11/08/12
Lab Sample ID:	D40798-1	Date Received:	11/10/12
Matrix:	SO - Soil	Percent Solids:	89.6
Project:	PCU 296-6A		

## General Chemistry

Analyte	Result	RL	Units	DF	Analyzed	By	Method
<b>prep: DEPT.OF AG, BOOK N9</b>							
Specific Conductivity	377	1.0	umhos/cm	1	11/14/12	JD	SM2510B-1997 MOD
Chromium, Hexavalent	< 1.0	1.0	mg/kg	1	11/12/12	KB	SW846 3060A/7196A
Chromium, Trivalent <sup>a</sup>	41.8	2.1	mg/kg	1	11/13/12 16:30	JM	SW846 3060/7196A M
Redox Potential Vs H2	141		mv	1	11/14/12	CT	ASTM D1498-76M
Solids, Percent	89.6		%	1	11/12/12	SWT	SM19 2540B M
pH	9.75		su	1	11/14/12 12:45	CT	SW846 9045D

(a) Calculated as: (Chromium) - (Chromium, Hexavalent)

RL = Reporting Limit

**Report of Analysis**

Page 1 of 1

<b>Client Sample ID:</b>	CUT 2 SUBLINER (COMP)	<b>Date Sampled:</b>	11/08/12
<b>Lab Sample ID:</b>	D40798-1A	<b>Date Received:</b>	11/10/12
<b>Matrix:</b>	SO - Soil	<b>Percent Solids:</b>	89.6
<b>Project:</b>	PCU 296-6A		

**SAR Metals Analysis**

Analyte	Result	RL	Units	DF	Prep	Analyzed By	Method	Prep Method
Calcium	15.3	2.0	mg/l	1	11/13/12	11/13/12 JM	SW846 6010C <sup>1</sup>	SW846 3010A/M <sup>2</sup>
Magnesium	4.56	1.0	mg/l	1	11/13/12	11/13/12 JM	SW846 6010C <sup>1</sup>	SW846 3010A/M <sup>2</sup>
Sodium	59.2	2.0	mg/l	1	11/13/12	11/13/12 JM	SW846 6010C <sup>1</sup>	SW846 3010A/M <sup>2</sup>

(1) Instrument QC Batch: MA2995

(2) Prep QC Batch: MP8877

RL = Reporting Limit

## Report of Analysis

Page 1 of 1

Client Sample ID:	CUT 2 SUBLINER (COMP)	Date Sampled:	11/08/12
Lab Sample ID:	D40798-1A	Date Received:	11/10/12
Matrix:	SO - Soil	Percent Solids:	89.6
Project:	PCU 296-6A		

## General Chemistry

Analyte	Result	RL	Units	DF	Analyzed	By	Method
Sodium Adsorption Ratio <sup>a</sup>	3.41		ratio	1	11/13/12 18:37	JM	USDA HANDBOOK 60

(a) Calculated as:  $(\text{Na meq/L}) / \sqrt{[(\text{Ca meq/L}) + (\text{Mg meq/L})/2]}$

---

RL = Reporting Limit



## Misc. Forms

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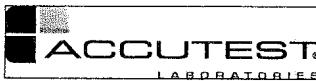
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### Custody Documents and Other Forms

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Includes the following where applicable:

- Chain of Custody



## **CHAIN OF CUSTODY**

PAGE 1 OF 1

Client / Reporting Information		Project Information		Requested Analysis (see TEST CODE sheet)		Matrix Codes	
Company Name <b>KRW Consulting</b>	Project Name: <b>XTO PCU 296-6A</b>	Street Address <b>8000 West 14th Street; Suite 200</b>	Street				
City <b>Lakewood, CO 80214</b>	City <b>XTO Energy</b>	State	Billing Information (if different from Report to)				
Project Contact <b>Dwayne Knudson</b>	Project # <b>1211-02</b>	Phone # <b>970-488-1098</b>	Client Purchase Order #	Street Address <b>21459 CR 5</b>	City <b>Rifle, CO 81650</b>		
Sampler(s) Name(s) <b>DAVID SANDERS</b>	Project Manager <b>Joe Hess</b>	Sampler(s) Name(s) <b>DAVID SANDERS</b>	Attention: <b>Jessica Dooling</b>				
Accutest Sample #	Field ID / Point of Collection <b>CRT 2 Submeter (COMP)</b>	Collection	Sampled by	Number of Preserved Bottles			
	MECH/DI Val #	Date <b>11-8-12</b>	Time <b>11:50</b>	Matrix <b>DS</b>	# of bottles <b>5</b>	ICP	
						NH4H	
						NH4O3	
						HSO4	
						NONE	
						D Water	
						MECH	
						ENCORE	
						Bisulfite	
						T - 910	
						X	
							<b>O1</b>
Turnaround Time (Business days)							
Approved By (Accutest PM): / Date:				Data Deliverable Information			
<input type="checkbox"/> Std. 10 Business Days <input checked="" type="checkbox"/> Std. 6 Business Days (By contract only) <input type="checkbox"/> 3 Day Emergency <input type="checkbox"/> 2 Day Emergency <input type="checkbox"/> 1 Day Emergency <input type="checkbox"/>				<input type="checkbox"/> Commercial "A" (Level 1) <input type="checkbox"/> Commercial "B" (Level 2) <input type="checkbox"/> COMMNB <input type="checkbox"/> COMMBN+			
				<input type="checkbox"/> State Forms Required <input type="checkbox"/> Send Forms to State <input type="checkbox"/> Report by Fax <input checked="" type="checkbox"/> Report by PDF ONLY <input type="checkbox"/> EDD Format			
				Please email to: <b>KRW Piceance Team</b>			
Commercial "A" = Results Only Commercial "B" = Results + QC Summary Commercial BN = Results/QC/Narrative (+ = chromatograms)							
Sample Custody must be documented below each time samples change possession, including courier delivery.							
Relinquished by Sampler: <b>1 Lori Abington</b>	Date Time: <b>11/9/12 15:30</b>	Received By: <b>1 White Service Center</b>	Relinquished By: <b>2</b>	Date Time:	Received By: <b>2 D J R 10/10</b>		
Relinquished by Sampler: <b>3</b>	Date Time:	Received By: <b>3</b>	Relinquished By: <b>4</b>	Date Time:	Received By: <b>4</b>		
Relinquished by: <b>5</b>	Date Time:	Received By: <b>5</b>	Custody Seal: <b>X</b>	Intact: <input checked="" type="checkbox"/>	Preserved where applicable	On Ice: <input checked="" type="checkbox"/>	Cooler Temp: <b>35</b>

## D40798: Chain of Custody

Page 1 of 2



## Accutest Laboratories Sample Receipt Summary

Accutest Job Number: D40798

Client: KRW CONSULTING

Immediate Client Services Action Required: No

Date / Time Received: 11/10/2012 10:15:00 A

No. Coolers:

1

Client Service Action Required at Login: No

Project: XTO PCU 296-6A

Airbill #'s: FX

### Cooler Security      Y or N

- |                           |                                     |                          |                       |                                     |                          |
|---------------------------|-------------------------------------|--------------------------|-----------------------|-------------------------------------|--------------------------|
| 1. Custody Seals Present: | <input checked="" type="checkbox"/> | <input type="checkbox"/> | 3. COC Present:       | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| 2. Custody Seals Intact:  | <input checked="" type="checkbox"/> | <input type="checkbox"/> | 4. Smpl Dates/Time OK | <input checked="" type="checkbox"/> | <input type="checkbox"/> |

### Cooler Temperature      Y or N

- |                              |                                     |                          |
|------------------------------|-------------------------------------|--------------------------|
| 1. Temp criteria achieved:   | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| 2. Cooler temp verification: | Infrared gun                        |                          |
| 3. Cooler media:             | Ice (bag)                           |                          |

### Quality Control Preservation      Y or N      N/A

- |                                 |                                     |                          |
|---------------------------------|-------------------------------------|--------------------------|
| 1. Trip Blank present / cooler: | <input type="checkbox"/>            | <input type="checkbox"/> |
| 2. Trip Blank listed on COC:    | <input type="checkbox"/>            | <input type="checkbox"/> |
| 3. Samples preserved properly:  | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| 4. VOCs headspace free:         | <input type="checkbox"/>            | <input type="checkbox"/> |

### Sample Integrity - Documentation

- |  |                                     |                          |
|--|-------------------------------------|--------------------------|
| 1. Sample labels present on bottles:   | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| 2. Container labeling complete:        | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| 3. Sample container label / COC agree: | <input checked="" type="checkbox"/> | <input type="checkbox"/> |

### Sample Integrity - Condition

- |                                  |                                     |                          |
|----------------------------------|-------------------------------------|--------------------------|
| 1. Sample recvd within HT:       | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| 2. All containers accounted for: | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| 3. Condition of sample:          | Intact                              |                          |

### Sample Integrity - Instructions

- |   |                                     |                                     |
|---|-------------------------------------|-------------------------------------|
| 1. Analysis requested is clear:           | <input checked="" type="checkbox"/> | <input type="checkbox"/>            |
| 2. Bottles received for unspecified tests | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |
| 3. Sufficient volume rec'd for analysis:  | <input checked="" type="checkbox"/> | <input type="checkbox"/>            |
| 4. Compositing instructions clear:        | <input type="checkbox"/>            | <input type="checkbox"/>            |
| 5. Filtering instructions clear:          | <input type="checkbox"/>            | <input checked="" type="checkbox"/> |

Comments

Accutest Laboratories  
V:(303) 425-6021

4036 Youngfield Street  
F: (303) 425-6854

Wheat Ridge, CO  
www.accutest.com

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D40798: Chain of Custody

Page 2 of 2



## GC/MS Volatiles

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### QC Data Summaries

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Includes the following where applicable:

- Method Blank Summaries
- Blank Spike Summaries
- Matrix Spike and Duplicate Summaries



**Method Blank Summary**

Job Number: D40798  
 Account: XTOKRWR XTO Energy  
 Project: PCU 296-6A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
V5V1501-MB	5V24597.D	1	11/14/12	BD	n/a	n/a	V5V1501

The QC reported here applies to the following samples:

Method: SW846 8260B

D40798-1

CAS No.	Compound	Result	RL	MDL	Units	Q
71-43-2	Benzene	ND	50	25	ug/kg	
100-41-4	Ethylbenzene	ND	100	19	ug/kg	
108-88-3	Toluene	ND	100	50	ug/kg	
1330-20-7	Xylene (total)	ND	200	100	ug/kg	

CAS No.	Surrogate Recoveries	Limits	
2037-26-5	Toluene-D8	103%	64-130%
460-00-4	4-Bromofluorobenzene	92%	62-131%
17060-07-0	1,2-Dichloroethane-D4	96%	70-130%

CAS No.	Surrogate Recoveries	Limits
2037-26-5	Toluene-D8	103%
460-00-4	4-Bromofluorobenzene	92%
17060-07-0	1,2-Dichloroethane-D4	96%

## Blank Spike Summary

Page 1 of 1

Job Number: D40798

Account: XTOKWR XTO Energy

Project: PCU 296-6A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
V5V1501-BS	5V24598.D	1	11/14/12	BD	n/a	n/a	V5V1501

The QC reported here applies to the following samples:

Method: SW846 8260B

D40798-1

CAS No.	Compound	Spike ug/kg	BSP ug/kg	BSP %	Limits
71-43-2	Benzene	50	43.6	87	70-130
100-41-4	Ethylbenzene	50	45.0	90	70-130
108-88-3	Toluene	50	44.2	88	70-130
1330-20-7	Xylene (total)	150	138	92	70-130

CAS No.	Surrogate Recoveries	BSP	Limits
2037-26-5	Toluene-D8	103%	64-130%
460-00-4	4-Bromofluorobenzene	97%	62-131%
17060-07-0	1,2-Dichloroethane-D4	100%	70-130%

\* = Outside of Control Limits.

# Matrix Spike/Matrix Spike Duplicate Summary

Page 1 of 1

Job Number: D40798

Account: XTOKWR XTO Energy

Project: PCU 296-6A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
D40713-1MS	5V24600.D	1	11/14/12	BD	n/a	n/a	V5V1501
D40713-1MSD	5V24601.D	1	11/14/12	BD	n/a	n/a	V5V1501
D40713-1	5V24599.D	1	11/14/12	BD	n/a	n/a	V5V1501

The QC reported here applies to the following samples:

Method: SW846 8260B

D40798-1

CAS No.	Compound	D40713-1		Spike	MS	MS	MSD	MSD	Limits	
		ug/kg	Q	ug/kg	ug/kg	%	ug/kg	%	RPD	Rec/RPD
71-43-2	Benzene	ND		3170	2820	89	2670	84	5	64-139/30
100-41-4	Ethylbenzene	ND		3170	2950	93	2740	86	7	68-136/30
108-88-3	Toluene	ND		3170	2870	91	2710	86	6	60-130/30
1330-20-7	Xylene (total)	ND		9510	9070	95	8450	89	7	58-142/30

CAS No.	Surrogate Recoveries	MS	MSD	D40713-1	Limits
2037-26-5	Toluene-D8	103%	102%	101%	64-130%
460-00-4	4-Bromofluorobenzene	105%	104%	99%	62-131%
17060-07-0	1,2-Dichloroethane-D4	97%	97%	99%	70-130%

\* = Outside of Control Limits.



## GC/MS Volatiles

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### Raw Data

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## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\V5111312.S\  
 Data File : 5V24609.D  
 Acq On : 14 Nov 2012 9:48 am  
 Operator : BRETD  
 Sample : D40798-1  
 Misc : MS4956,V5V1501,5.019,,100,5,1  
 ALS Vial : 38 Sample Multiplier: 1

Quant Time: Nov 14 16:50:58 2012  
 Quant Method : C:\msdchem\1\METHODS\V5AP1497TVH1497.M  
 Quant Title : 8260  
 QLast Update : Wed Nov 14 09:56:27 2012  
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
2) Pentafluorobenzene	11.613	168	423379	50.00	ug/l	-0.01
35) 1,4-Difluorobenzene	12.412	114	512936	50.00	ug/l	-0.01
53) Chlorobenzene-d5	15.072	117	463620	50.00	ug/l	0.00
74) 1,4-Dichlorobenzene-d4	17.036	152	347427	50.00	ug/l	0.00

System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	12.012	102	32533	45.15	ug/l	-0.01
Spiked Amount 50.000	Range 70 - 130		Recovery	=	90.30%	
61) Toluene-d8	13.816	98	563347	51.29	ug/l	0.00
Spiked Amount 50.000	Range 70 - 130		Recovery	=	102.58%	
69) 4-Bromofluorobenzene	16.008	95	238114	50.30	ug/l	-0.01
Spiked Amount 50.000	Range 70 - 130		Recovery	=	100.60%	

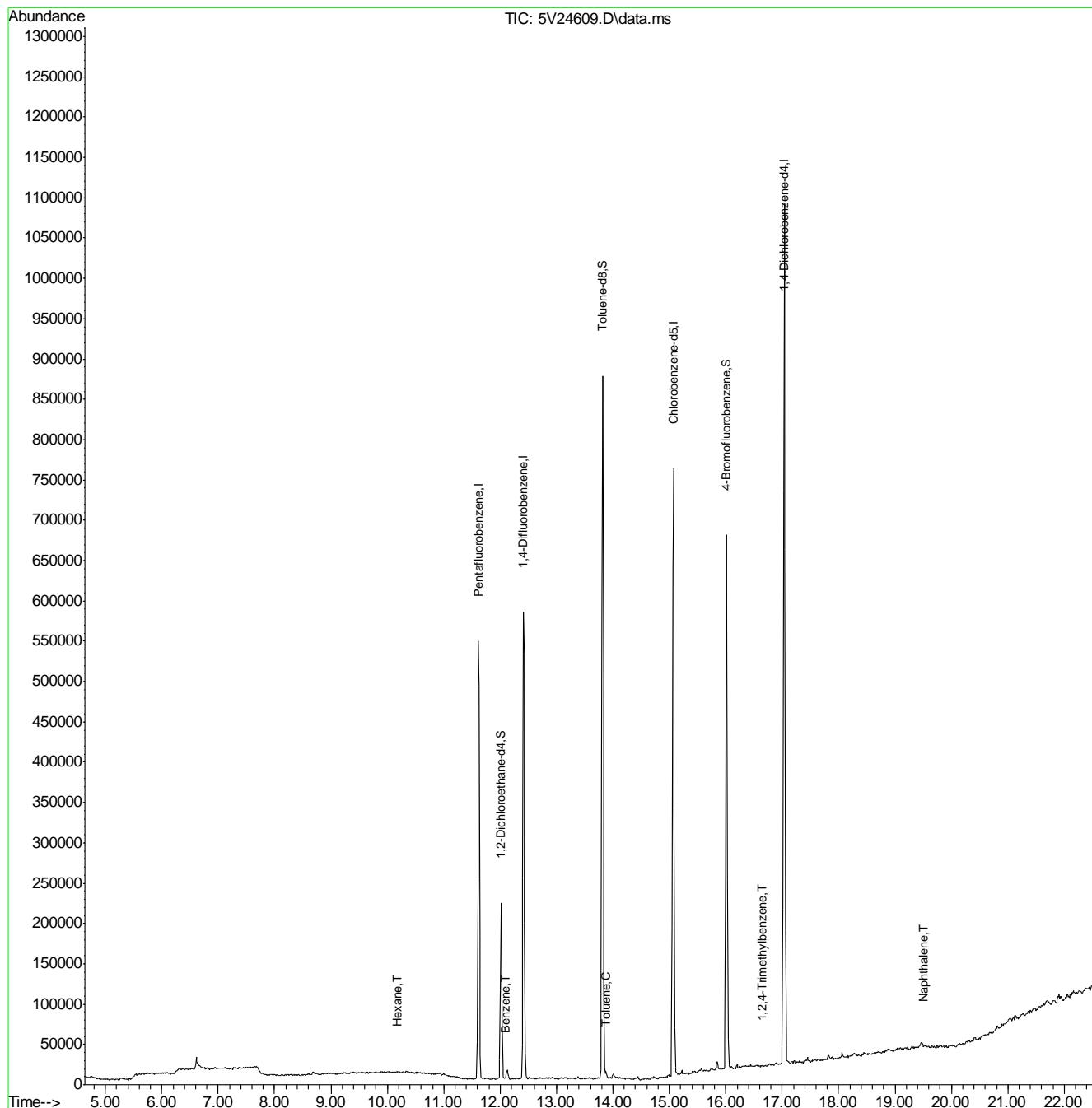
Target Compounds					QValue
41) Hexane	10.174	57	193	0.04	ug/l 100
50) Benzene	12.092	78	1830	0.13	ug/l 100
62) Toluene	13.873	92	2715	0.32	ug/l # 85
82) 1,2,4-Trimethylbenzene	16.648	105	949	0.06	ug/l # 83
91) Naphthalene	19.513	128	1034	0.06	ug/l 100

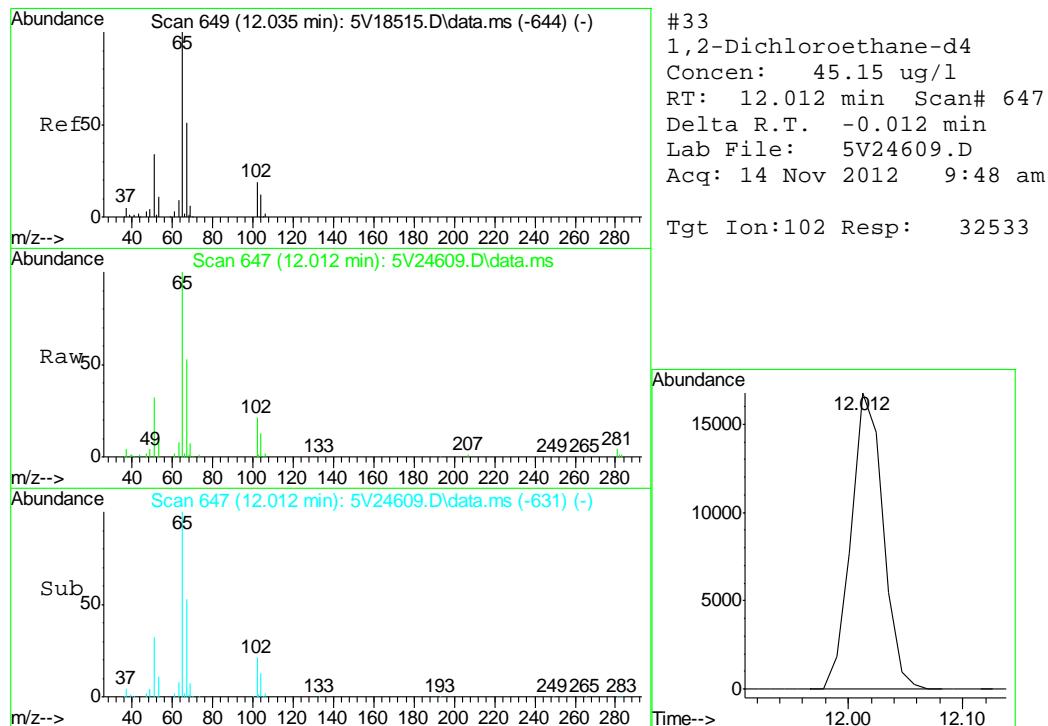
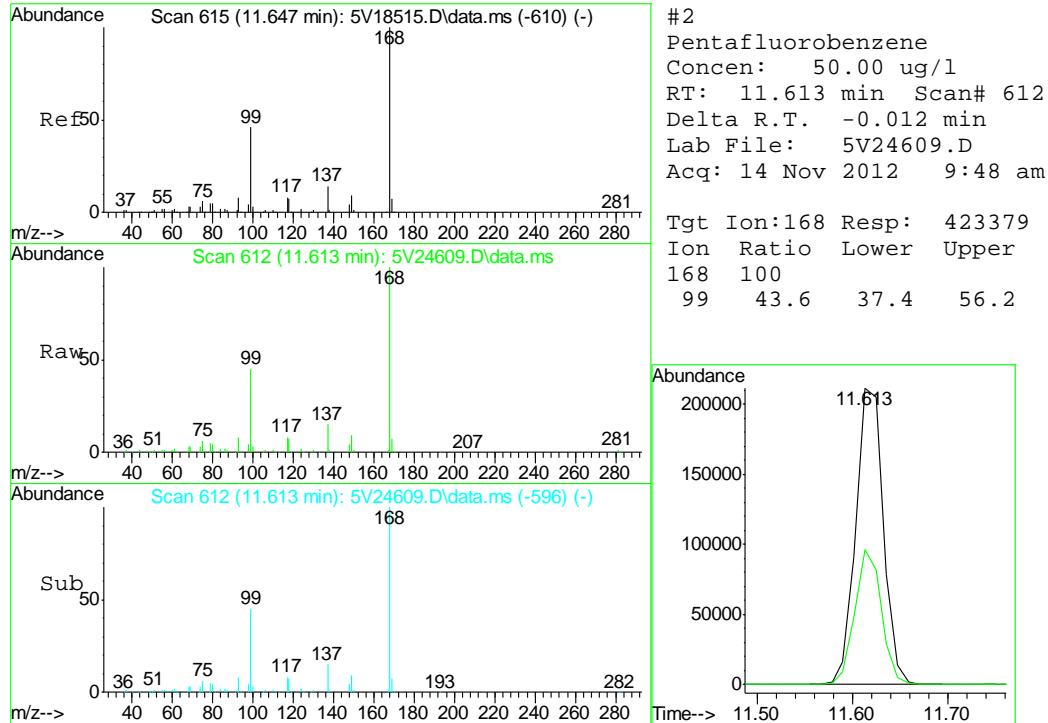
(#) = qualifier out of range (m) = manual integration (+) = signals summed

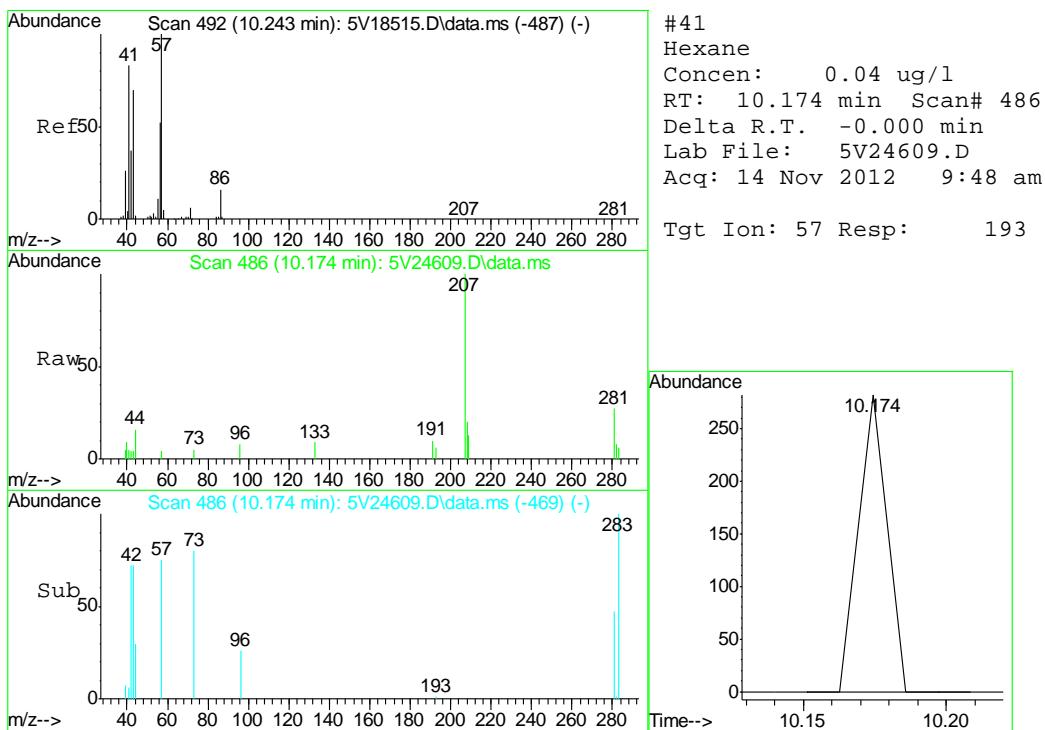
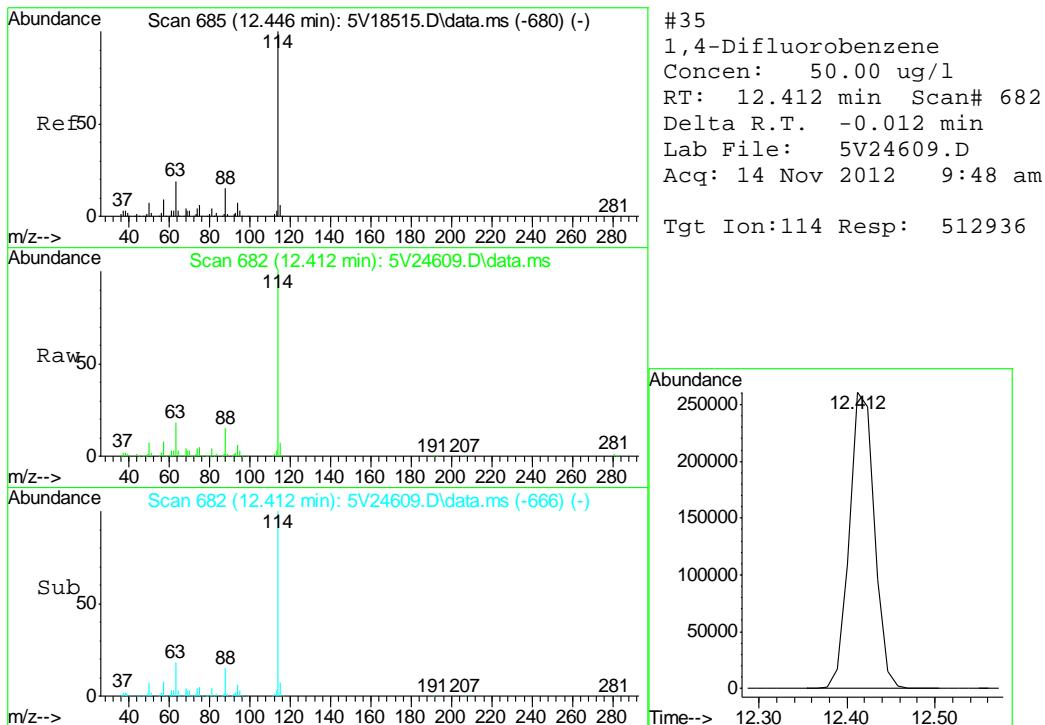
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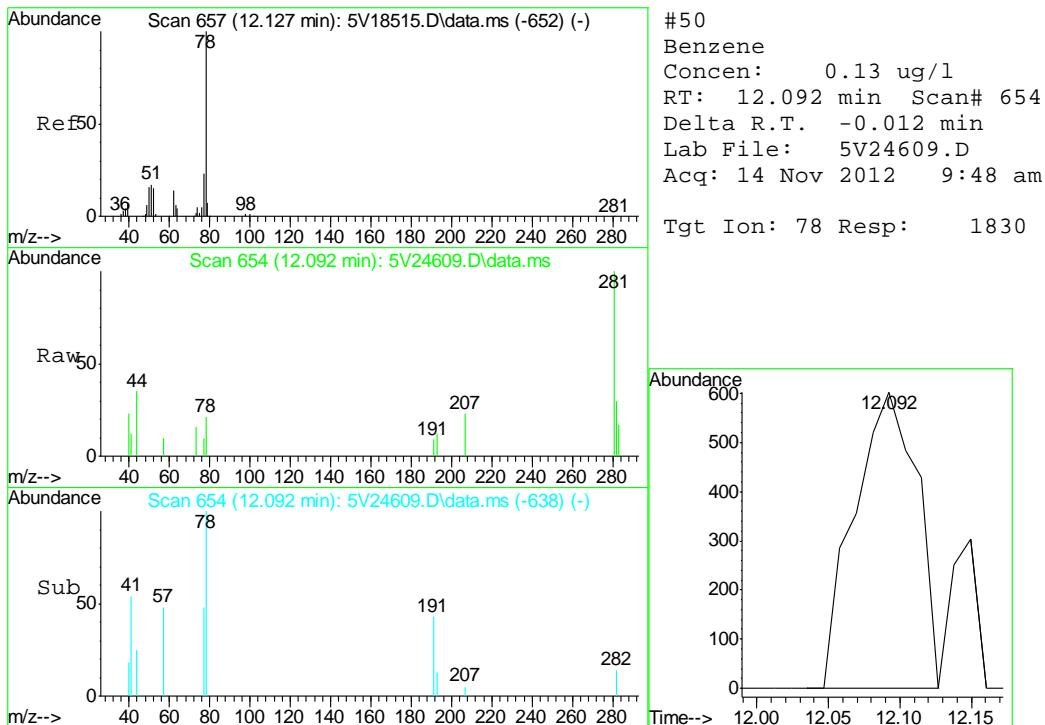
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 Data File : 5V24609.D  
 Acq On : 14 Nov 2012 9:48 am  
 Operator : BRETD  
 Sample : D40798-1  
 Misc : MS4956,V5V1501,5.019,,100,5,1  
 ALS Vial : 38 Sample Multiplier: 1

Quant Time: Nov 14 16:50:58 2012  
 Quant Method : C:\msdchem\1\METHODS\V5AP1497TVH1497.M  
 Quant Title : 8260  
 QLast Update : Wed Nov 14 09:56:27 2012  
 Response via : Initial Calibration

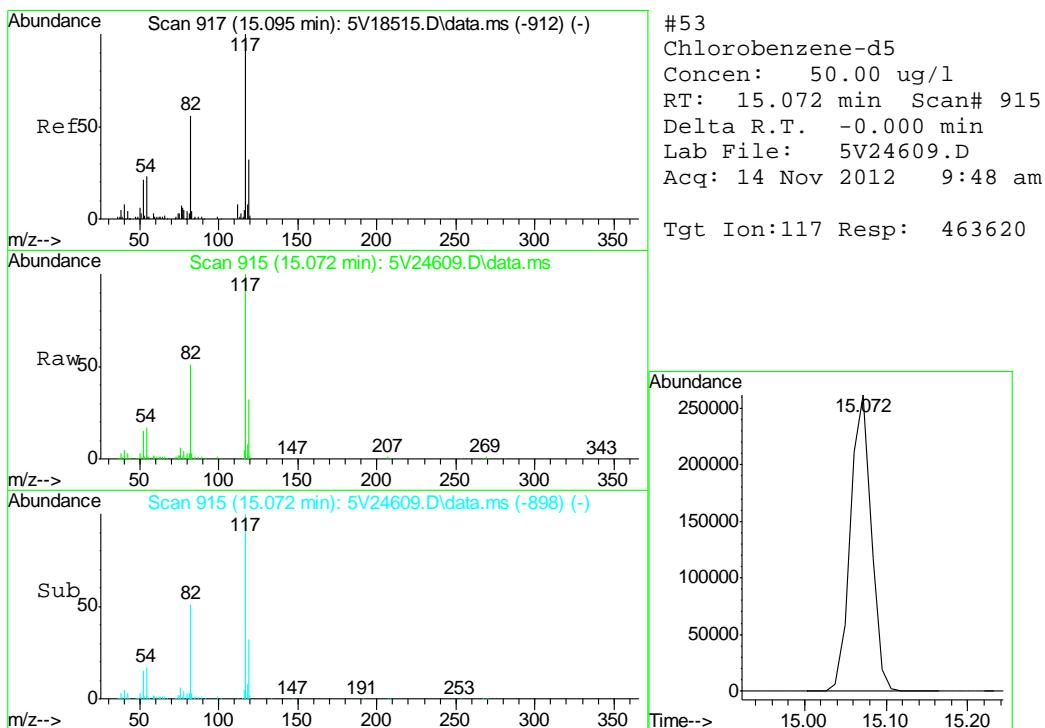


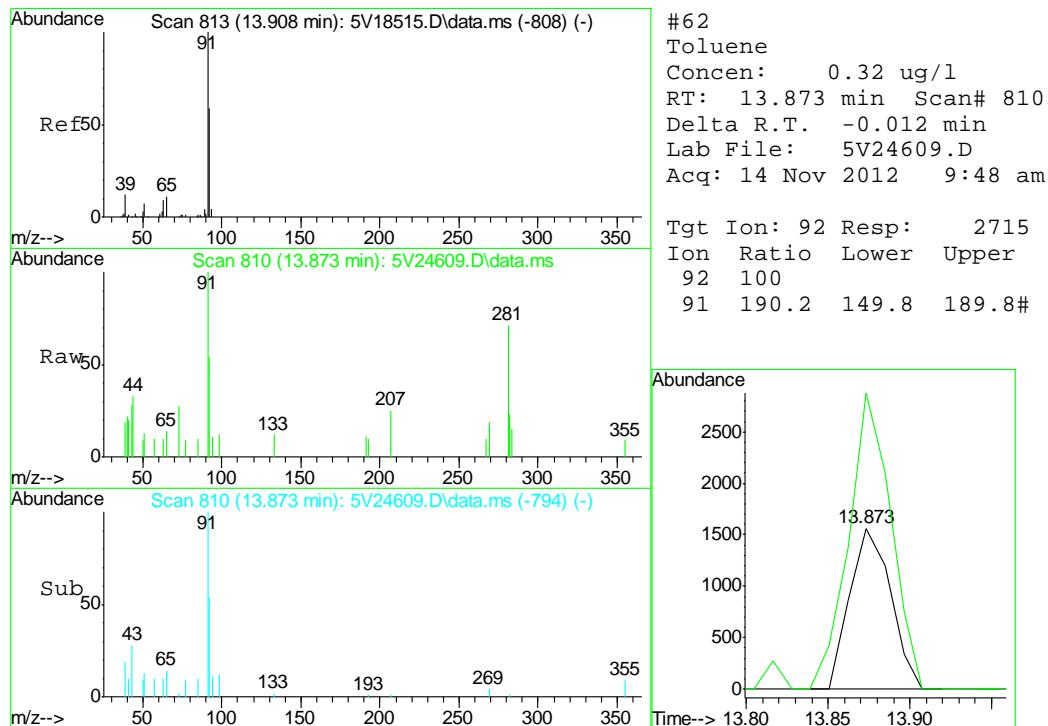
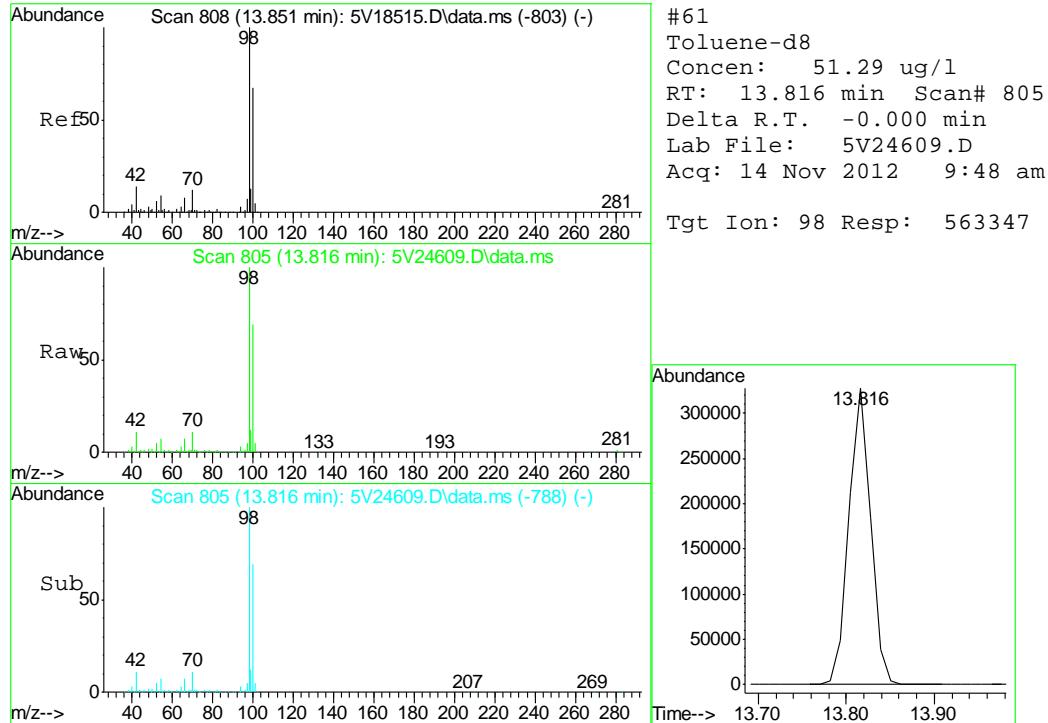


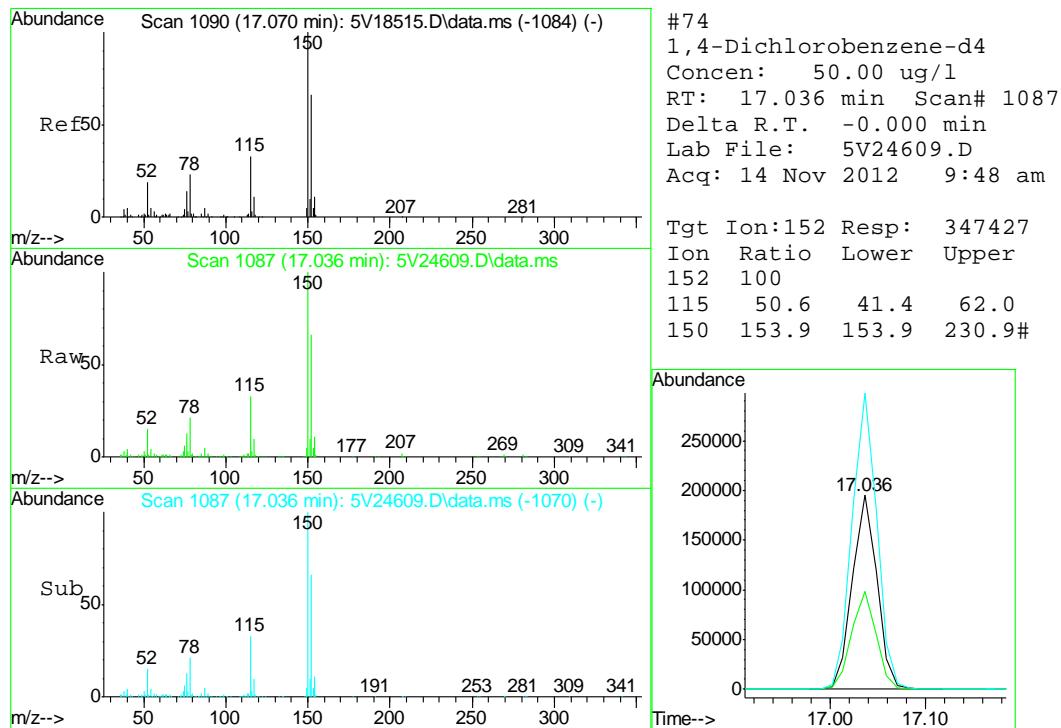
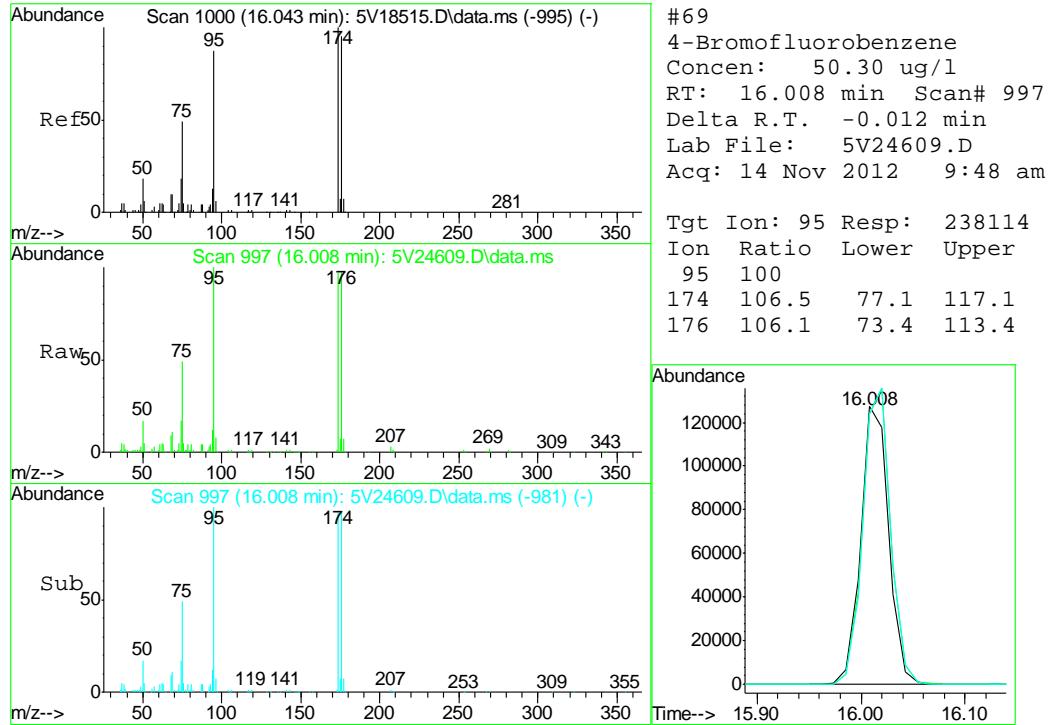


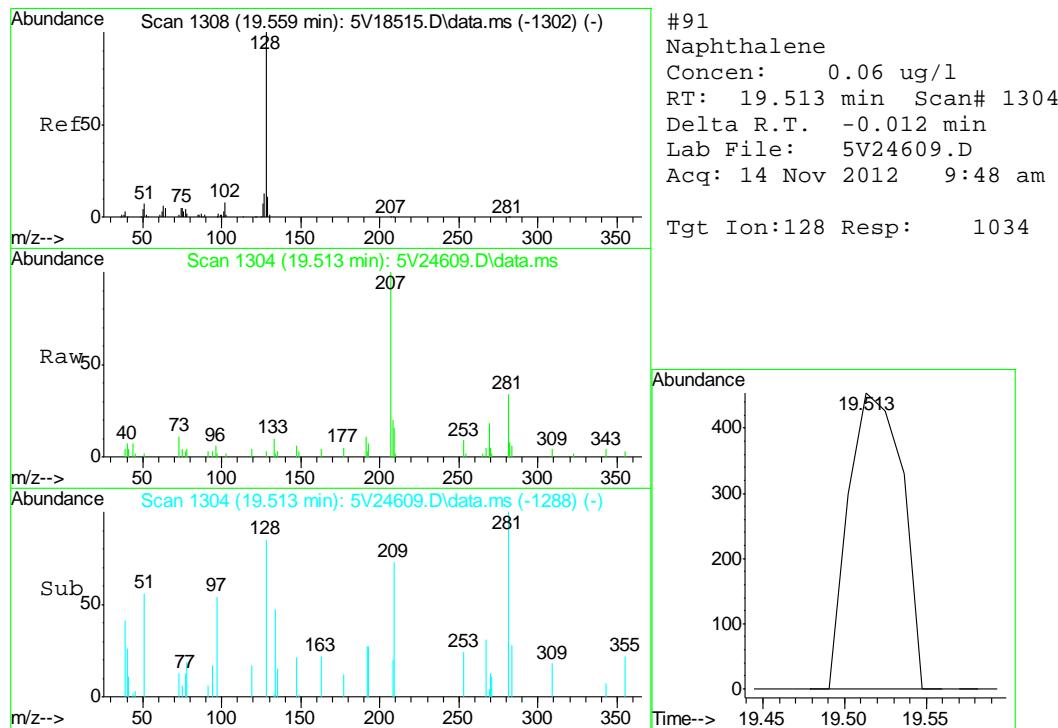
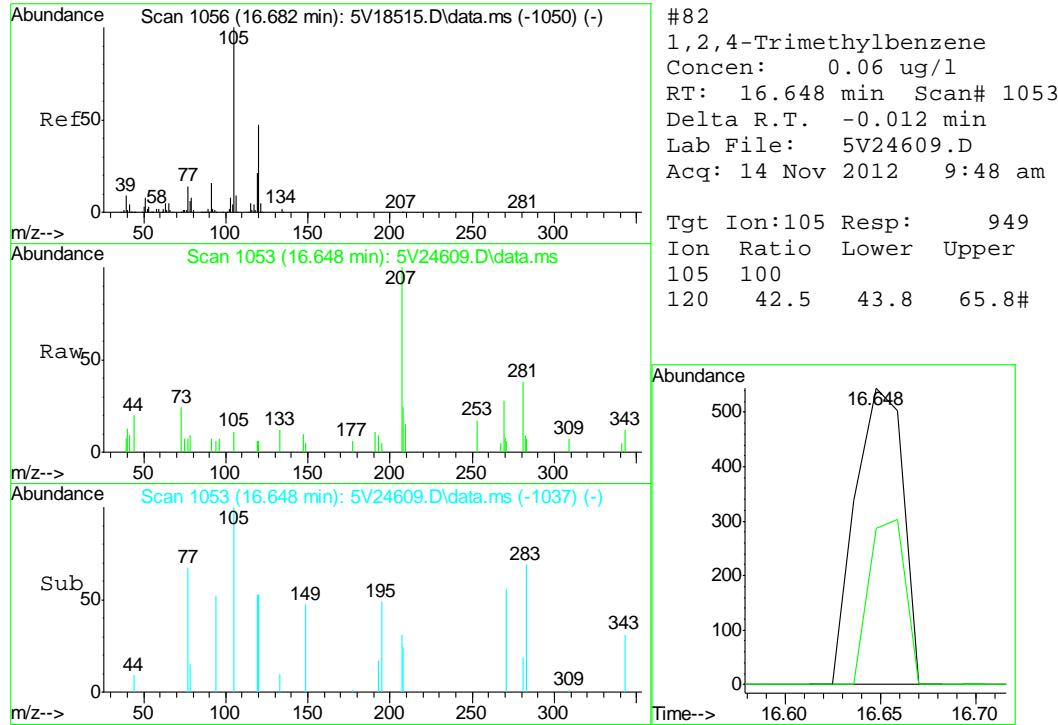


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## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\V5111312.S\  
 Data File : 5V24597.D  
 Acq On : 14 Nov 2012 3:00 am  
 Operator : BRETD  
 Sample : MB  
 Misc : MS4956,V5V1501,5.00,,100,5,1  
 ALS Vial : 26 Sample Multiplier: 1

Quant Time: Nov 14 15:14:20 2012  
 Quant Method : C:\msdchem\1\METHODS\V5AP1497TVH1497.M  
 Quant Title : 8260  
 QLast Update : Wed Nov 14 09:56:27 2012  
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
2) Pentafluorobenzene	11.625	168	466596	50.00	ug/l	0.00
35) 1,4-Difluorobenzene	12.424	114	566385	50.00	ug/l	0.00
53) Chlorobenzene-d5	15.072	117	516924	50.00	ug/l	0.00
74) 1,4-Dichlorobenzene-d4	17.036	152	338090	50.00	ug/l	0.00

System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	12.024	102	38234	48.14	ug/l	0.00
Spiked Amount 50.000	Range 70 - 130		Recovery	=	96.28%	
61) Toluene-d8	13.817	98	630619	51.50	ug/l	0.00
Spiked Amount 50.000	Range 70 - 130		Recovery	=	103.00%	
69) 4-Bromofluorobenzene	16.020	95	243711	46.17	ug/l	0.00
Spiked Amount 50.000	Range 70 - 130		Recovery	=	92.34%	

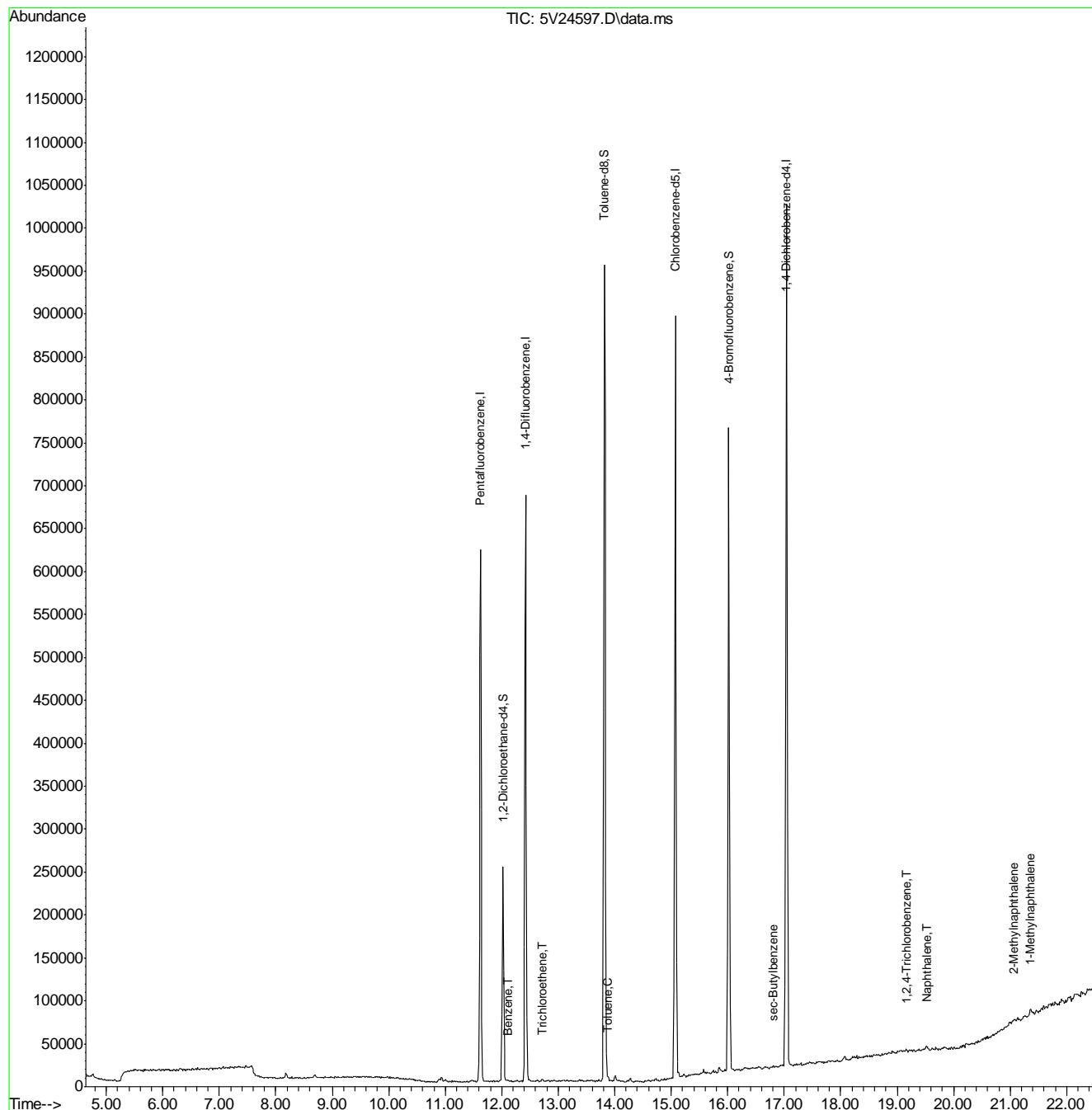
Target Compounds					Qvalue
1) TVH-Gasoline	13.102	TIC	141616m	Below Cal	
48) Trichloroethene	12.721	95	565	0.14 ug/l	88
50) Benzene	12.104	78	1340	0.09 ug/l	100
62) Toluene	13.874	92	854	0.09 ug/l #	63
83) sec-Butylbenzene	16.808	105	1026	0.05 ug/l	97
90) 1,2,4-Trichlorobenzene	19.160	180	1663	0.18 ug/l #	79
91) Naphthalene	19.525	128	5416	0.32 ug/l	100
94) 2-Methylnaphthalene	21.066	142	2952	5.30 ug/l #	67
95) 1-Methylnaphthalene	21.363	142	3770	1.39 ug/l #	94

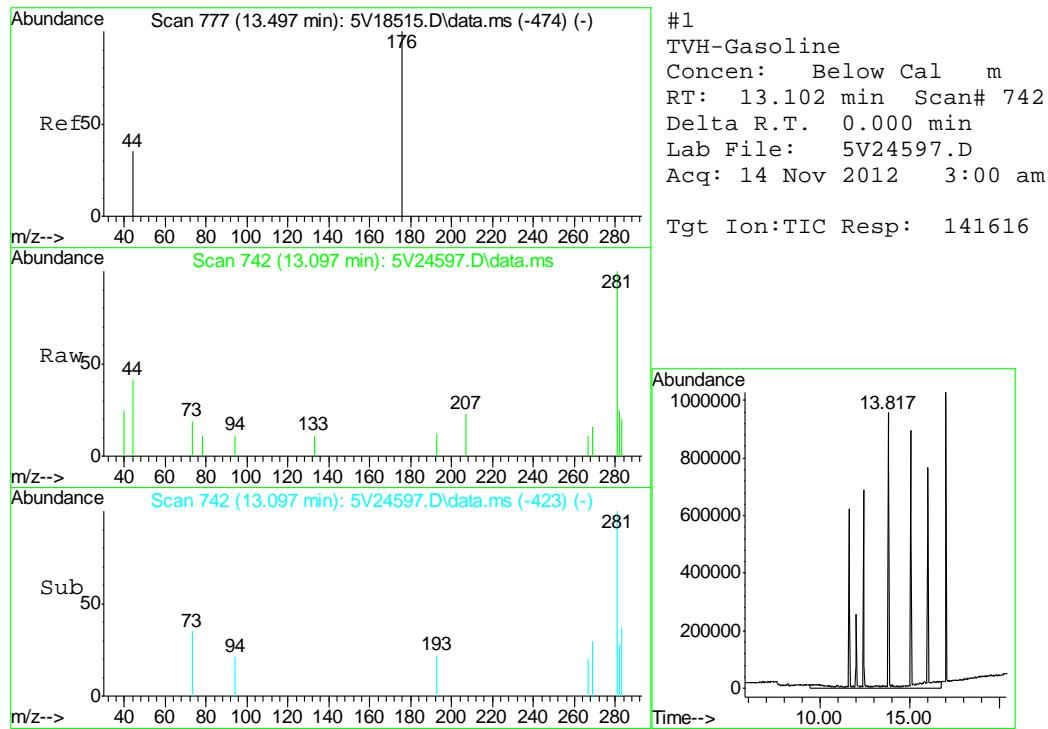
(#) = qualifier out of range (m) = manual integration (+) = signals summed

## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\V5111312.S\  
 Data File : 5V24597.D  
 Acq On : 14 Nov 2012 3:00 am  
 Operator : BRETD  
 Sample : MB  
 Misc : MS4956,V5V1501,5.00,,100,5,1  
 ALS Vial : 26 Sample Multiplier: 1

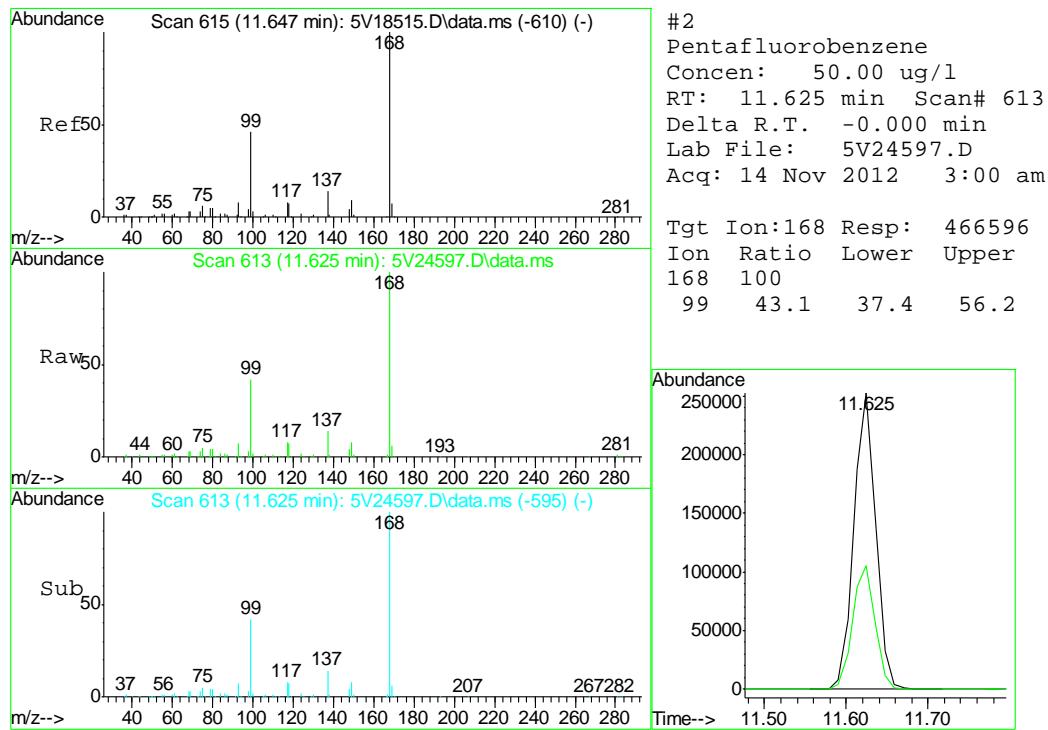
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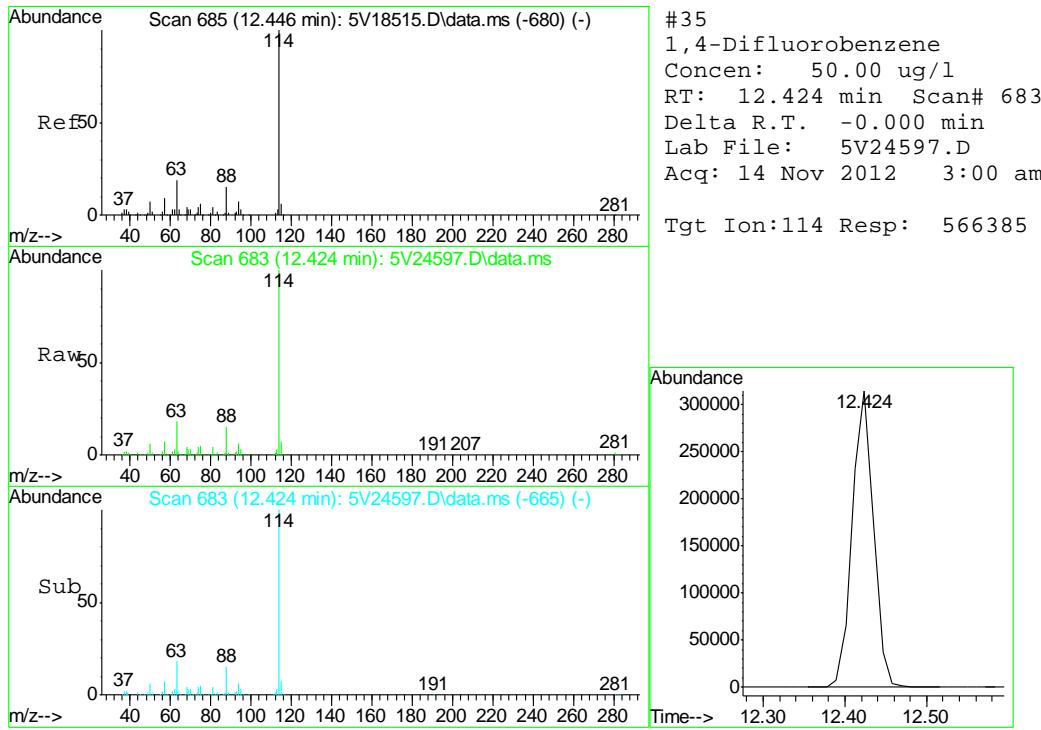
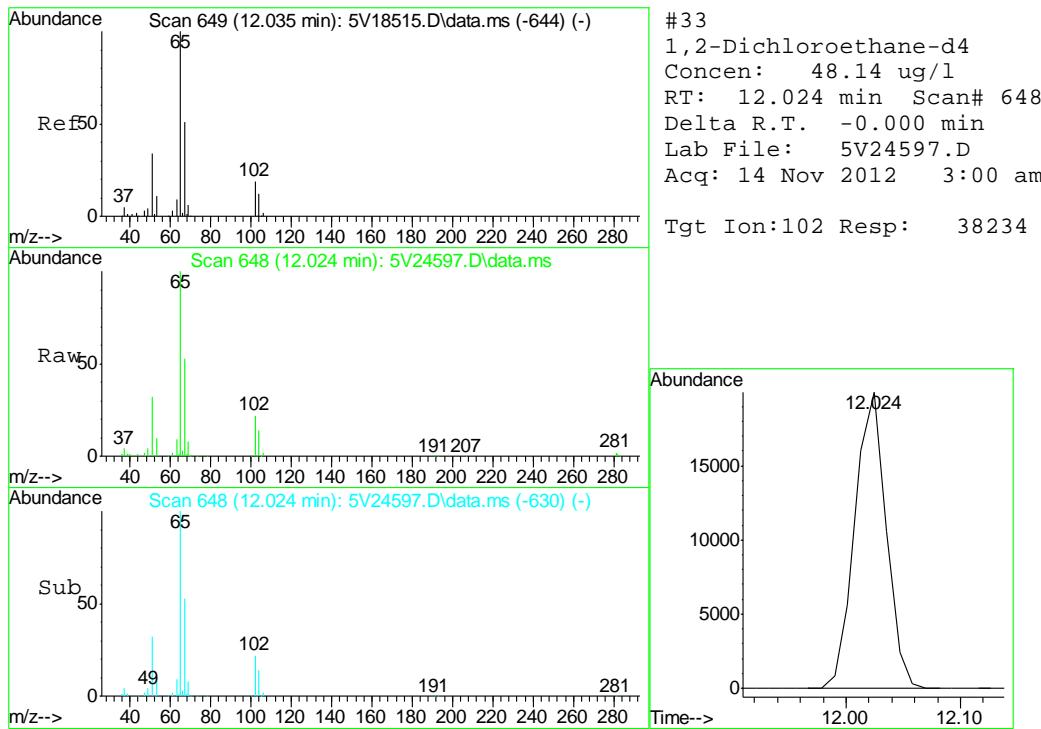


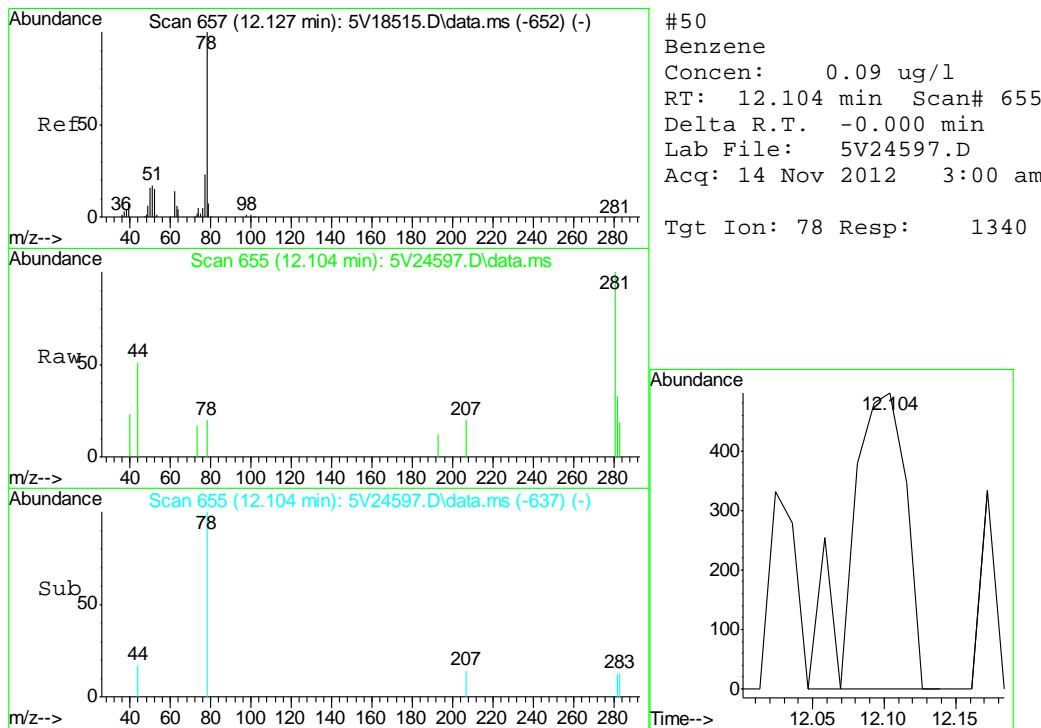
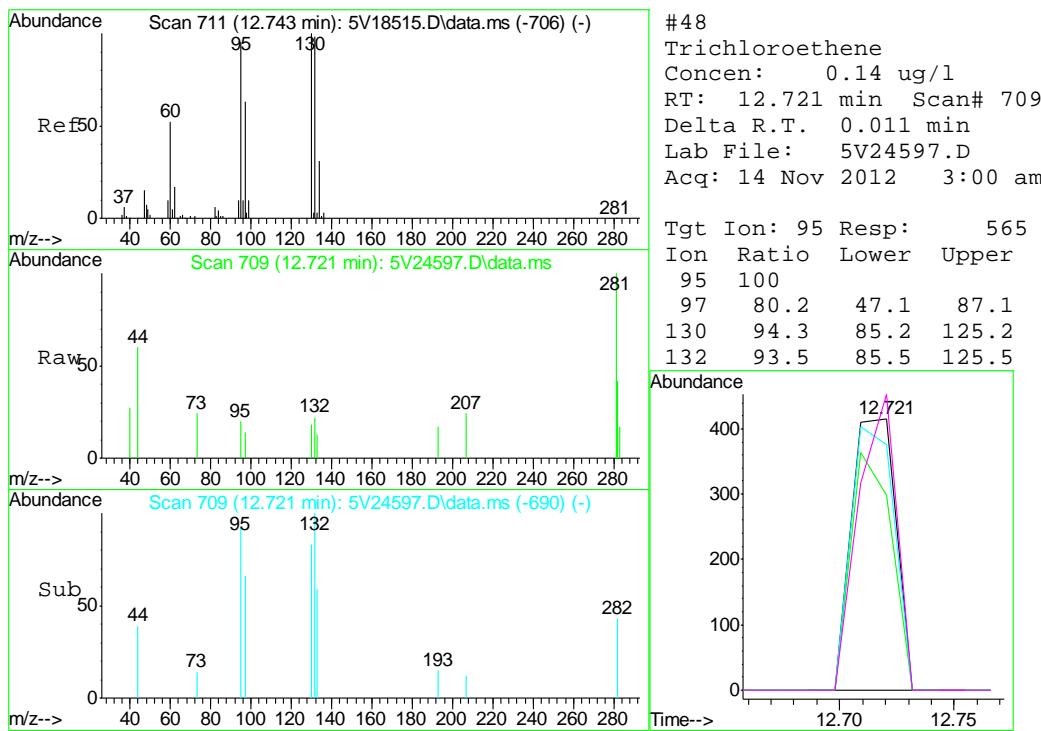


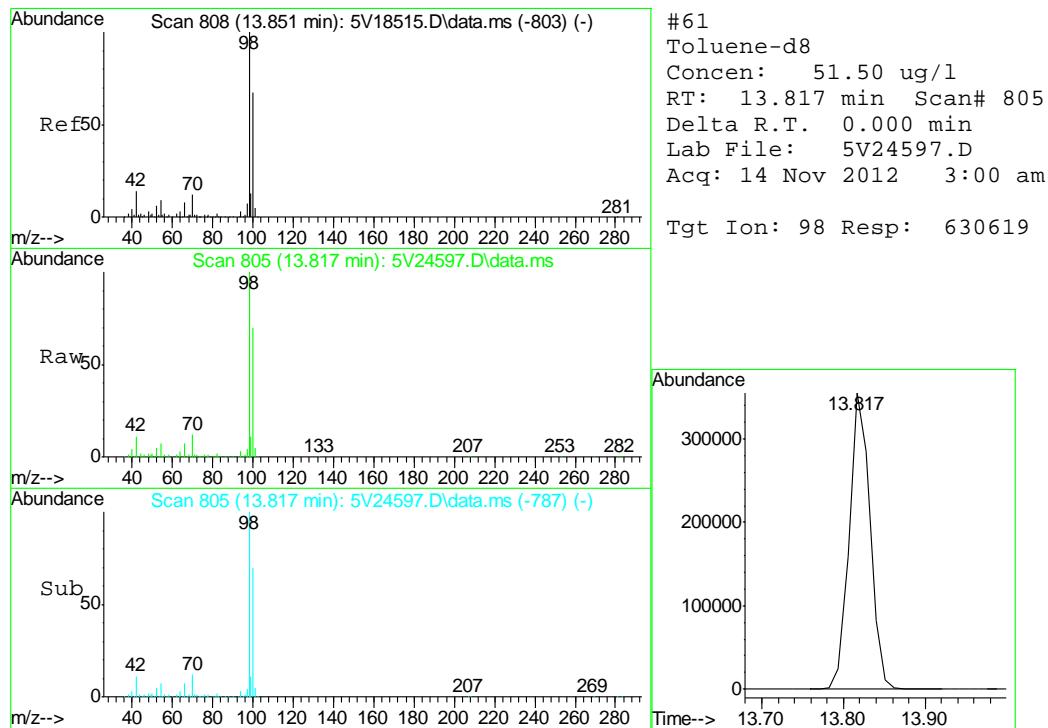
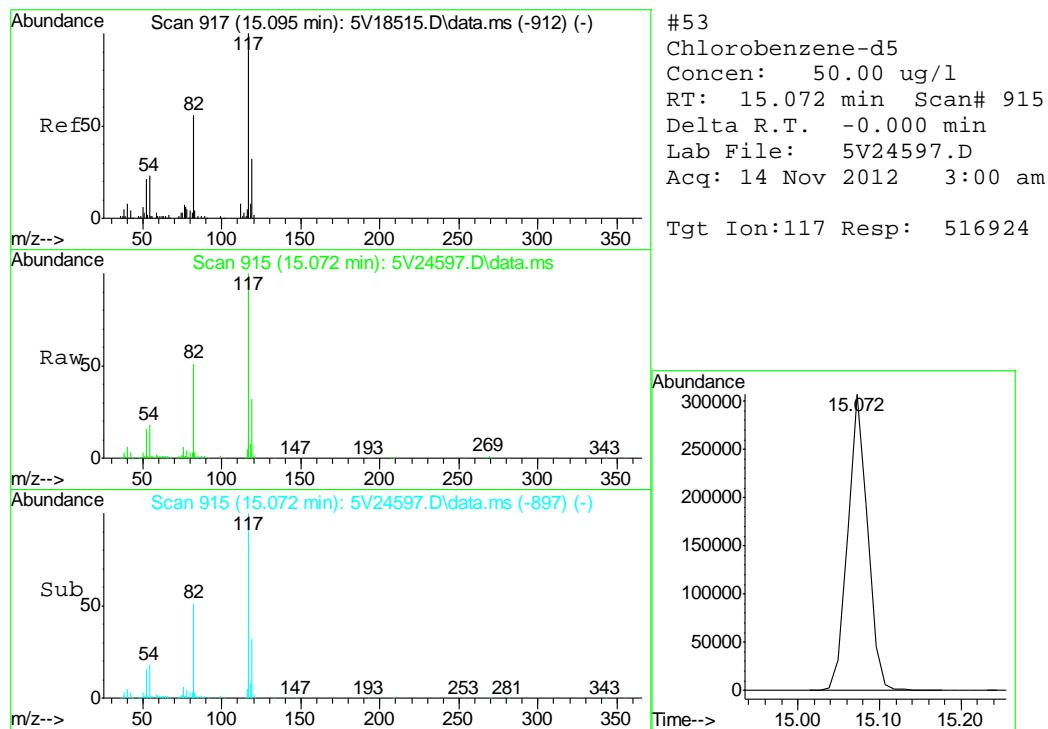
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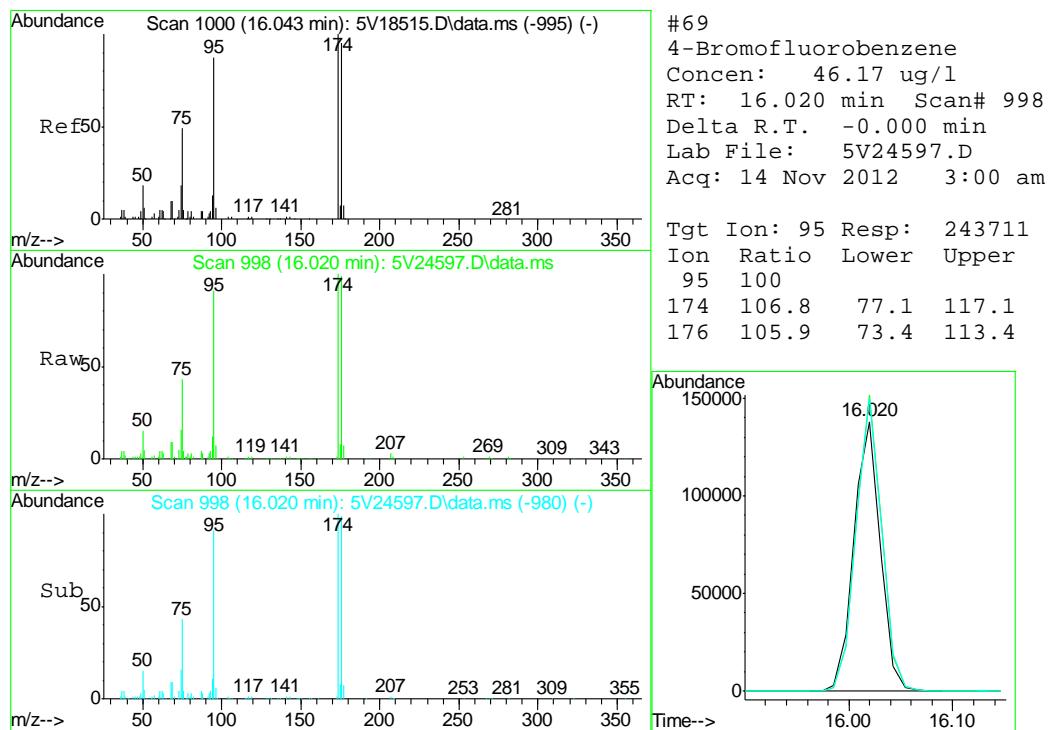
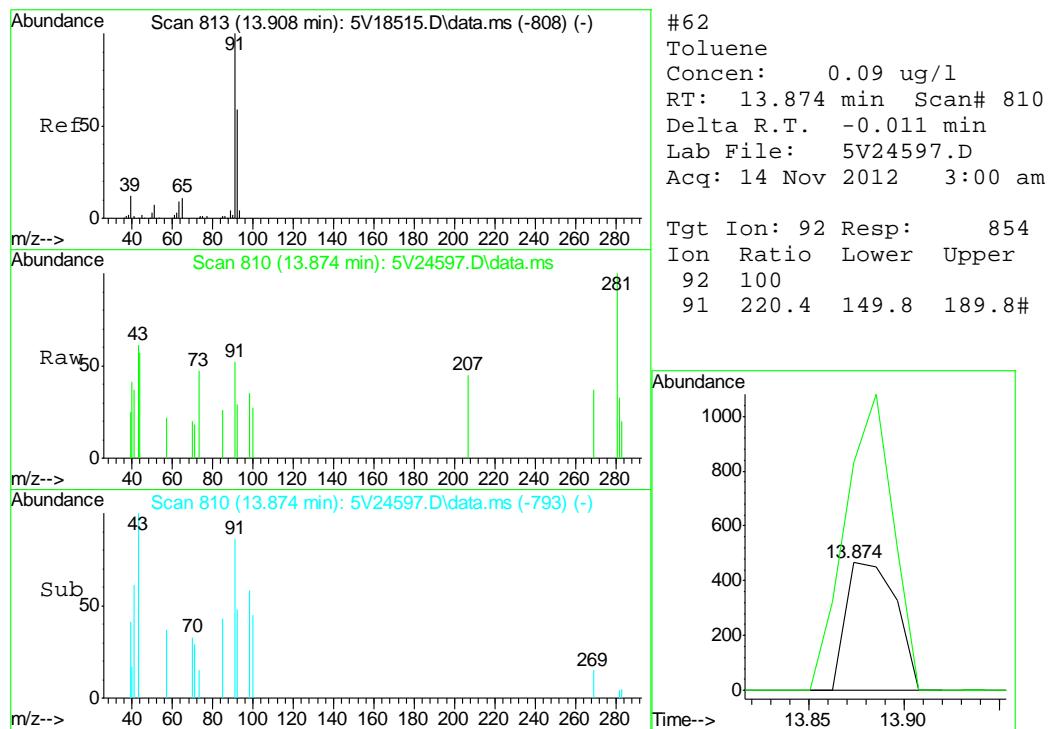
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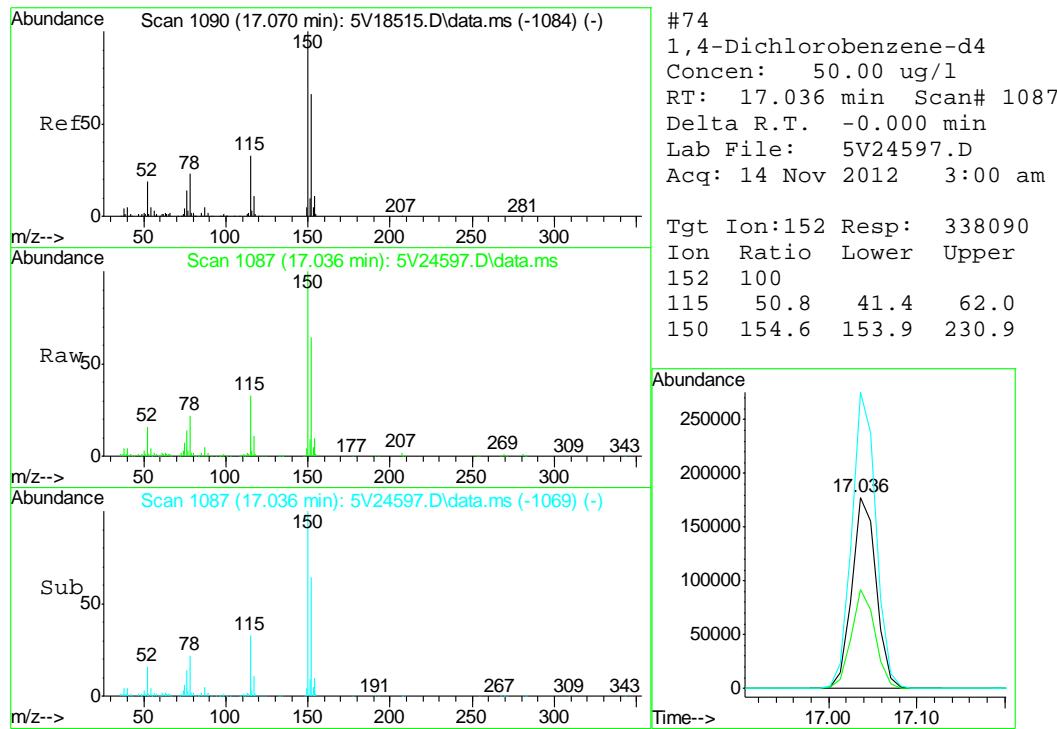






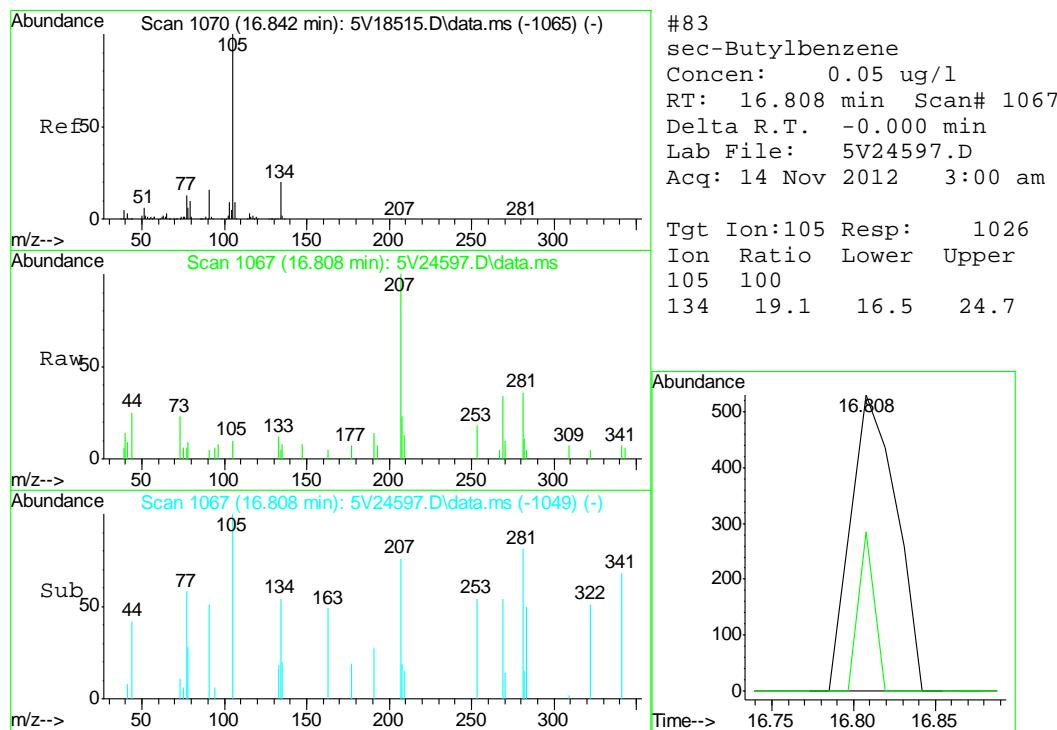


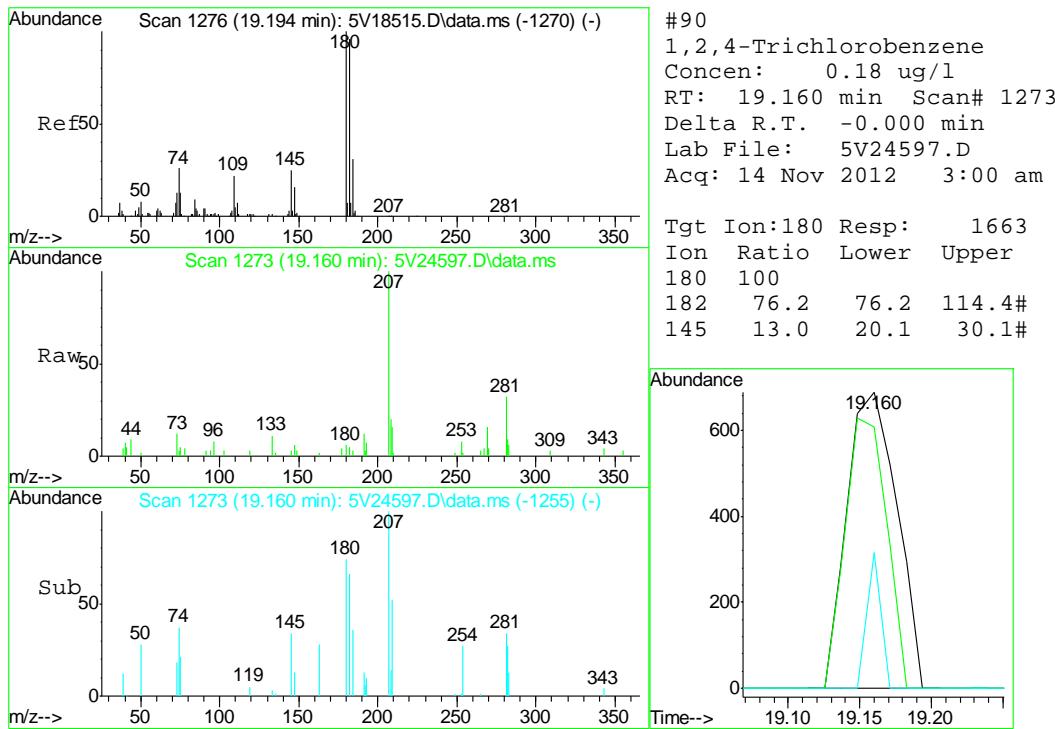




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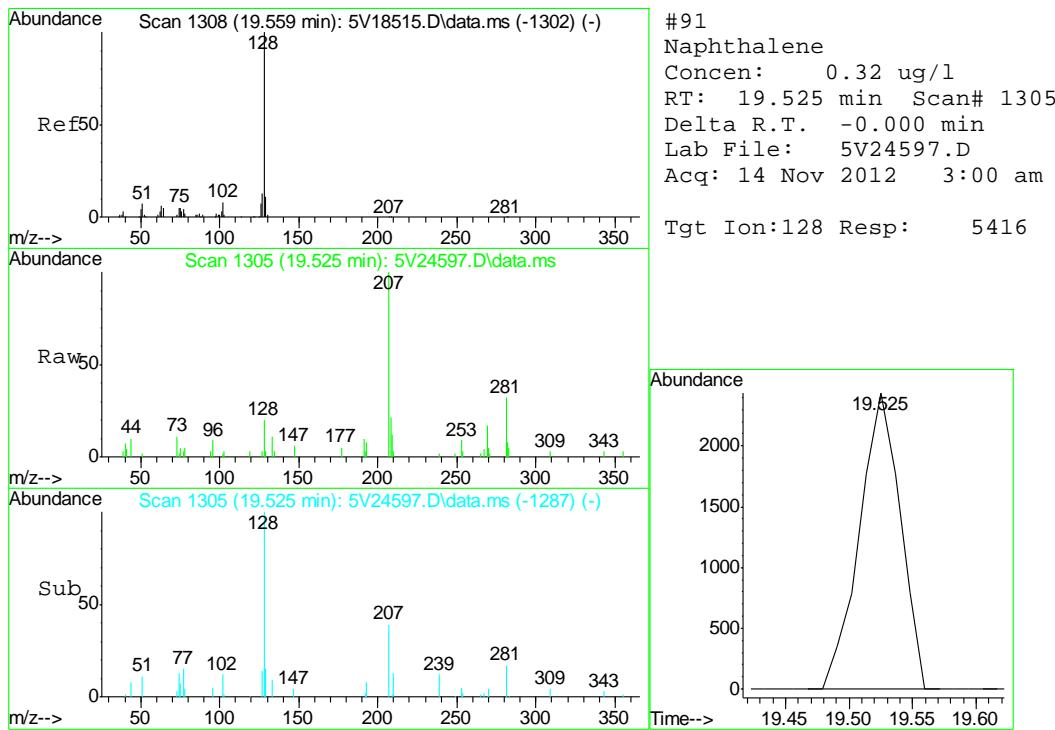
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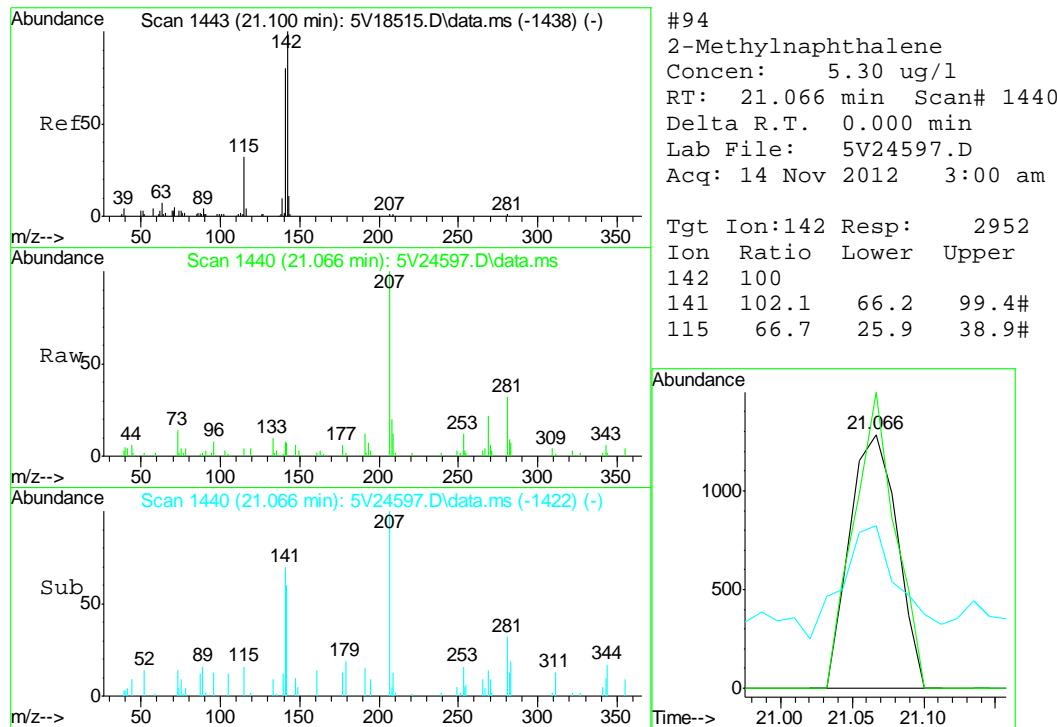




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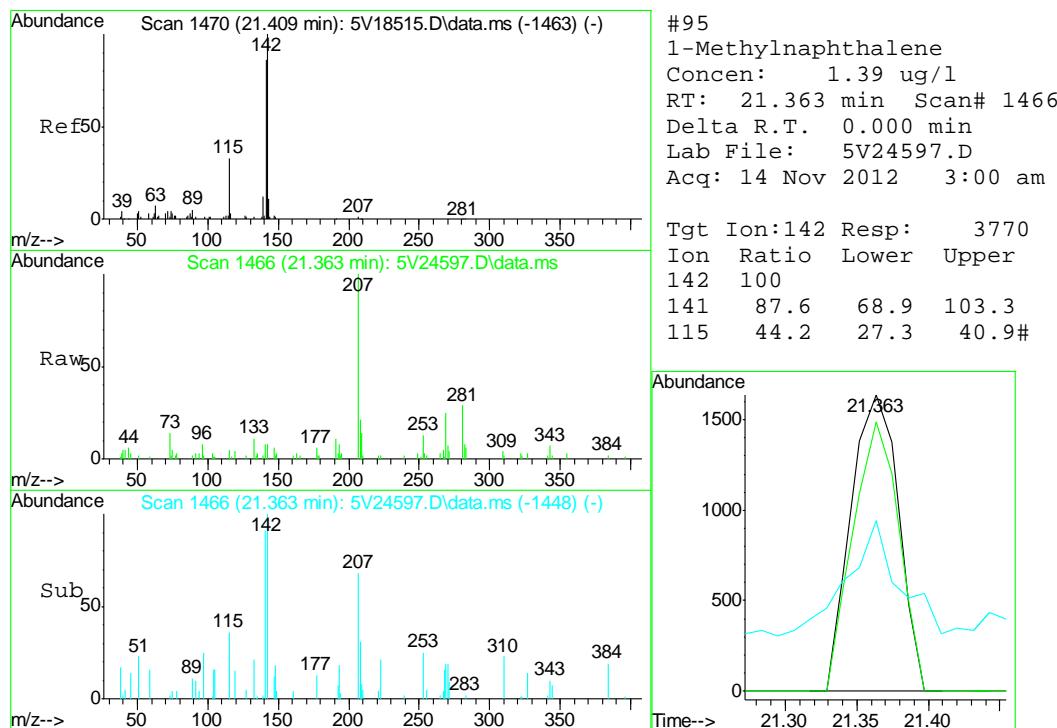
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## GC/MS Semi-volatiles

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### QC Data Summaries

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Includes the following where applicable:

- Method Blank Summaries
- Blank Spike Summaries
- Matrix Spike and Duplicate Summaries

**Method Blank Summary**

Job Number: D40798  
 Account: XTOKRWR XTO Energy  
 Project: PCU 296-6A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
OP6947-MB	3G12026.D	1	11/12/12	DC	11/12/12	OP6947	E3G568

The QC reported here applies to the following samples:

Method: SW846 8270C BY SIM

D40798-1

CAS No.	Compound	Result	RL	MDL	Units	Q
83-32-9	Acenaphthene	ND	8.3	4.3	ug/kg	
120-12-7	Anthracene	ND	8.3	4.3	ug/kg	
56-55-3	Benzo(a)anthracene	ND	8.3	4.3	ug/kg	
205-99-2	Benzo(b)fluoranthene	ND	8.3	4.3	ug/kg	
207-08-9	Benzo(k)fluoranthene	ND	8.3	4.3	ug/kg	
50-32-8	Benzo(a)pyrene	ND	8.3	4.3	ug/kg	
218-01-9	Chrysene	ND	8.3	4.3	ug/kg	
53-70-3	Dibenzo(a,h)anthracene	ND	8.3	4.3	ug/kg	
206-44-0	Fluoranthene	ND	8.3	4.3	ug/kg	
86-73-7	Fluorene	ND	8.3	4.3	ug/kg	
193-39-5	Indeno(1,2,3-cd)pyrene	ND	8.3	4.3	ug/kg	
91-20-3	Naphthalene	ND	12	10	ug/kg	
129-00-0	Pyrene	ND	8.3	4.3	ug/kg	

CAS No.	Surrogate Recoveries	Limits
4165-60-0	Nitrobenzene-d5	84%
321-60-8	2-Fluorobiphenyl	80%
1718-51-0	Terphenyl-d14	94%

## Blank Spike Summary

Page 1 of 1

Job Number: D40798  
Account: XTOKRWR XTO Energy  
Project: PCU 296-6A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
OP6947-BS	3G12027.D	1	11/12/12	DC	11/12/12	OP6947	E3G568

The QC reported here applies to the following samples:

Method: SW846 8270C BY SIM

D40798-1

CAS No.	Compound	Spike ug/kg	BSP ug/kg	BSP %	Limits
83-32-9	Acenaphthene	83.3	79.5	95	68-130
120-12-7	Anthracene	83.3	80.2	96	67-130
56-55-3	Benzo(a)anthracene	83.3	79.0	95	65-130
205-99-2	Benzo(b)fluoranthene	83.3	84.4	101	44-130
207-08-9	Benzo(k)fluoranthene	83.3	80.5	97	56-131
50-32-8	Benzo(a)pyrene	83.3	82.1	99	62-130
218-01-9	Chrysene	83.3	81.6	98	70-130
53-70-3	Dibenzo(a,h)anthracene	83.3	81.5	98	55-130
206-44-0	Fluoranthene	83.3	76.5	92	70-130
86-73-7	Fluorene	83.3	75.7	91	70-130
193-39-5	Indeno(1,2,3-cd)pyrene	83.3	82.6	99	56-130
91-20-3	Naphthalene	83.3	78.0	94	70-130
129-00-0	Pyrene	83.3	84.5	101	70-130

CAS No.	Surrogate Recoveries	BSP	Limits
4165-60-0	Nitrobenzene-d5	81%	10-159%
321-60-8	2-Fluorobiphenyl	80%	19-131%
1718-51-0	Terphenyl-d14	89%	18-150%

\* = Outside of Control Limits.

# Matrix Spike/Matrix Spike Duplicate Summary

Page 1 of 1

Job Number: D40798

Account: XTOKWR XTO Energy

Project: PCU 296-6A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
OP6947-MS	3G12030.D	1	11/12/12	DC	11/12/12	OP6947	E3G568
OP6947-MSD	3G12031.D	1	11/12/12	DC	11/12/12	OP6947	E3G568
D40799-1	3G12029.D	1	11/12/12	DC	11/12/12	OP6947	E3G568

The QC reported here applies to the following samples:

Method: SW846 8270C BY SIM

D40798-1

CAS No.	Compound	D40799-1 ug/kg	Spike ug/kg	MS ug/kg	MS %	MSD ug/kg	MSD %	RPD	Limits Rec/RPD
83-32-9	Acenaphthene	ND	90.3	81.0	90	78.4	87	3	25-151/30
120-12-7	Anthracene	ND	90.3	89.0	99	89.4	99	0	39-159/30
56-55-3	Benzo(a)anthracene	ND	90.3	88.4	98	88.1	98	0	39-168/30
205-99-2	Benzo(b)fluoranthene	ND	90.3	97.5	108	94.3	104	3	24-163/30
207-08-9	Benzo(k)fluoranthene	ND	90.3	82.9	92	80.0	89	4	10-188/30
50-32-8	Benzo(a)pyrene	ND	90.3	84.6	94	84.3	93	0	32-144/30
218-01-9	Chrysene	ND	90.3	89.9	100	89.8	99	0	43-150/30
53-70-3	Dibenzo(a,h)anthracene	ND	90.3	86.8	96	86.4	96	0	21-152/30
206-44-0	Fluoranthene	ND	90.3	87.3	97	87.5	97	0	36-157/30
86-73-7	Fluorene	ND	90.3	83.3	92	79.9	88	4	10-182/30
193-39-5	Indeno(1,2,3-cd)pyrene	ND	90.3	87.8	97	87.1	96	1	20-154/30
91-20-3	Naphthalene	ND	90.3	82.9	92	78.8	87	5	10-163/30
129-00-0	Pyrene	ND	90.3	97.0	107	95.9	106	1	25-180/30

CAS No.	Surrogate Recoveries	MS	MSD	D40799-1	Limits
4165-60-0	Nitrobenzene-d5	71%	69%	73%	10-159%
321-60-8	2-Fluorobiphenyl	65%	64%	65%	19-131%
1718-51-0	Terphenyl-d14	79%	79%	75%	18-150%

\* = Outside of Control Limits.

8.3.1  
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## GC/MS Semi-volatiles

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### Raw Data

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## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\111212\  
 Data File : 3g12033.D  
 Acq On : 12 Nov 2012 7:00 pm  
 Operator : DONC  
 Sample : D40798-1  
 Misc : OP6947,E3G568,30.09,,,1,1  
 ALS Vial : 19 Sample Multiplier: 1

Quant Time: Nov 13 09:53:39 2012  
 Quant Method : C:\msdchem\1\METHODS\SIMPE3G568.M  
 Quant Title : PAHSIM BASE  
 QLast Update : Mon Nov 12 15:44:05 2012  
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Naphthalene-d8	5.789	136	146515	4.0000	ug/mL	0.00
6) Acenaphthene-d10	7.507	164	96104	4.0000	ug/mL	0.00
15) Phenanthrene-d10	8.995	188	154063	4.0000	ug/mL	0.00
19) Chrysene-d12	11.630	240	114035	4.0000	ug/mL	0.00
24) Perylene-d12	13.035	264	64180	4.0000	ug/mL	-0.01

System Monitoring Compounds						
2) Nitrobenzene-d5	5.103	82	539258	36.9495	ug/mL	0.00
Spiked Amount	50.000	Range	25 - 135	Recovery	=	73.90%
7) 2-Fluorobiphenyl	6.846	172	1483834	32.6327	ug/mL	0.00
Spiked Amount	50.000	Range	25 - 135	Recovery	=	65.26%
21) Terphenyl-d14	10.586	244	624848	39.4652	ug/mL	0.00
Spiked Amount	50.000	Range	25 - 135	Recovery	=	78.94%

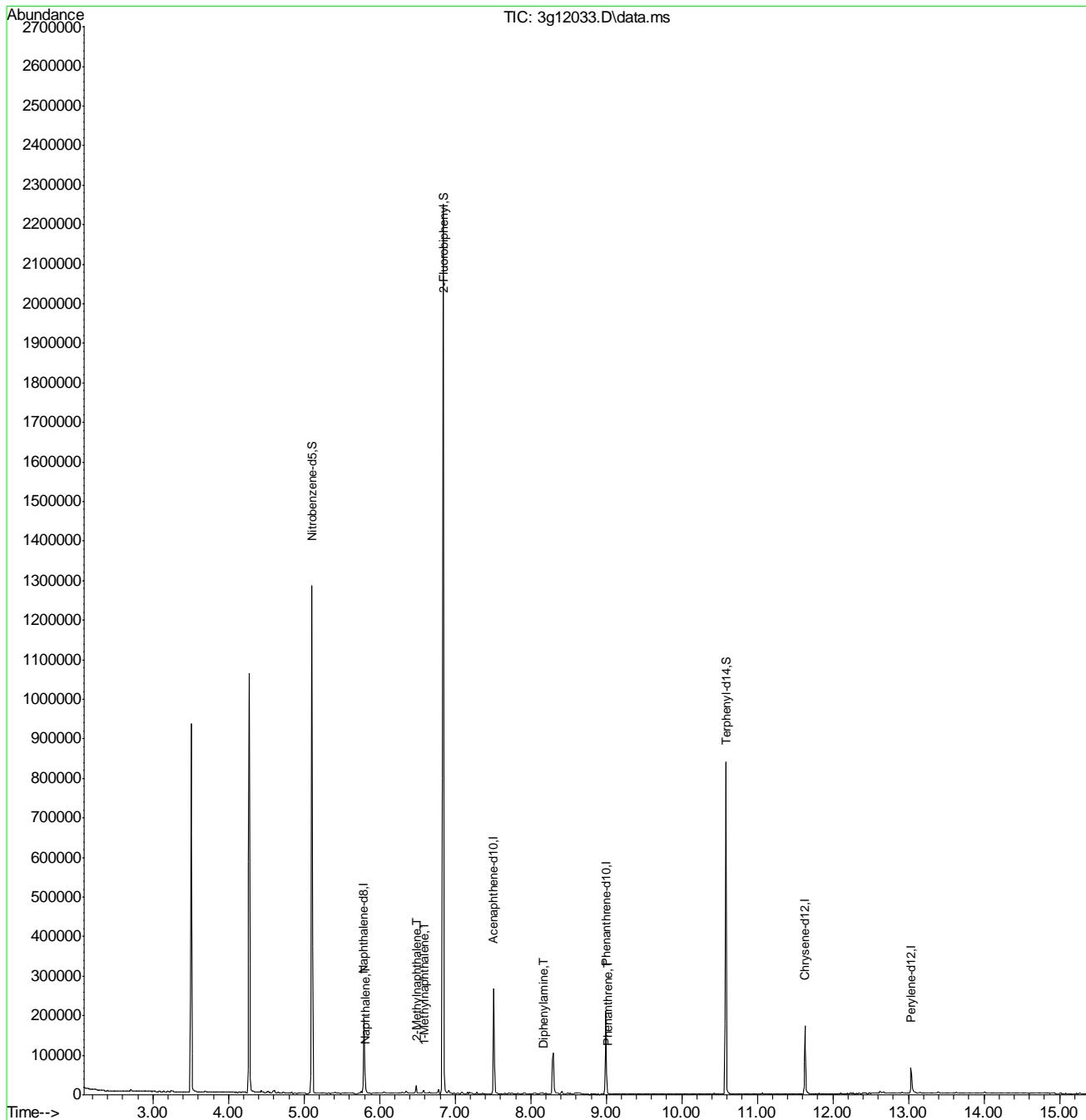
Target Compounds					Qvalue
3) N-Nitrosodimethylamine	2.487	74	11	N.D.	
4) N-Nitrosodi-propylamine	0.000	70	0	N.D. d	
5) Naphthalene	5.814	128	6077	0.1230 ug/mL	94
8) 2-Methylnaphthalene	6.487	142	8225	0.2324 ug/mL	98
9) 1-Methylnaphthalene	6.587	142	3672	0.1118 ug/mL	92
10) Acenaphthylene	7.366	152	530	N.D.	
11) Acenaphthene	7.507	154	469	Below Cal #	59
12) Dibenzofuran	7.720	168	1046	N.D.	
13) Fluorene	8.063	166	1185	N.D.	
14) Diphenylamine	8.169	169	2186	0.0611 ug/mL	72
16) Phenanthrene	9.019	178	3844	0.0670 ug/mL#	62
17) Anthracene	9.067	178	444	N.D.	
18) Fluoranthene	10.198	202	1098	N.D.	
20) Pyrene	10.428	202	1403	N.D.	
22) Benzo(a)anthracene	11.623	228	1258	N.D.	
23) Chrysene	11.623	228	1258	N.D.	
25) Benzo(b)fluoranthene	12.646	252	1670	N.D.	
26) Benzo(k)fluoranthene	12.646	252	1670	N.D.	
27) Benzo(a)pyrene	12.982	252	454	N.D.	
28) Indeno(1,2,3-cd)pyrene	14.318	276	560	N.D.	
29) Dibenz(a,h)anthracene	14.329	278	247	N.D.	
30) Benzo(g,h,i)perylene	14.318	276	586	N.D.	

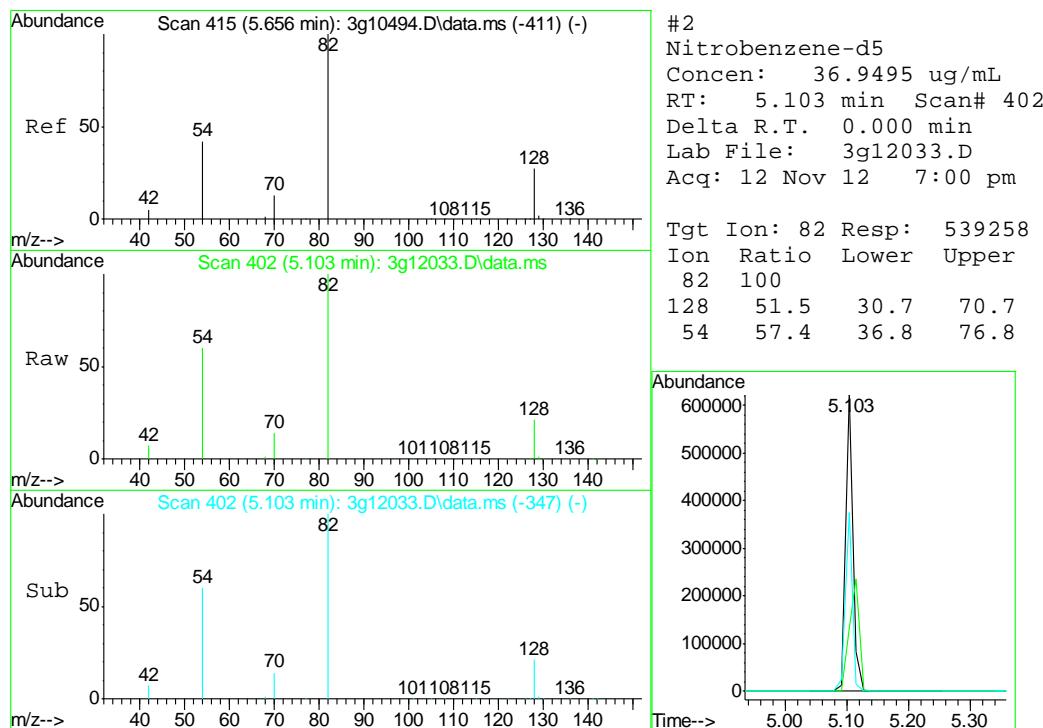
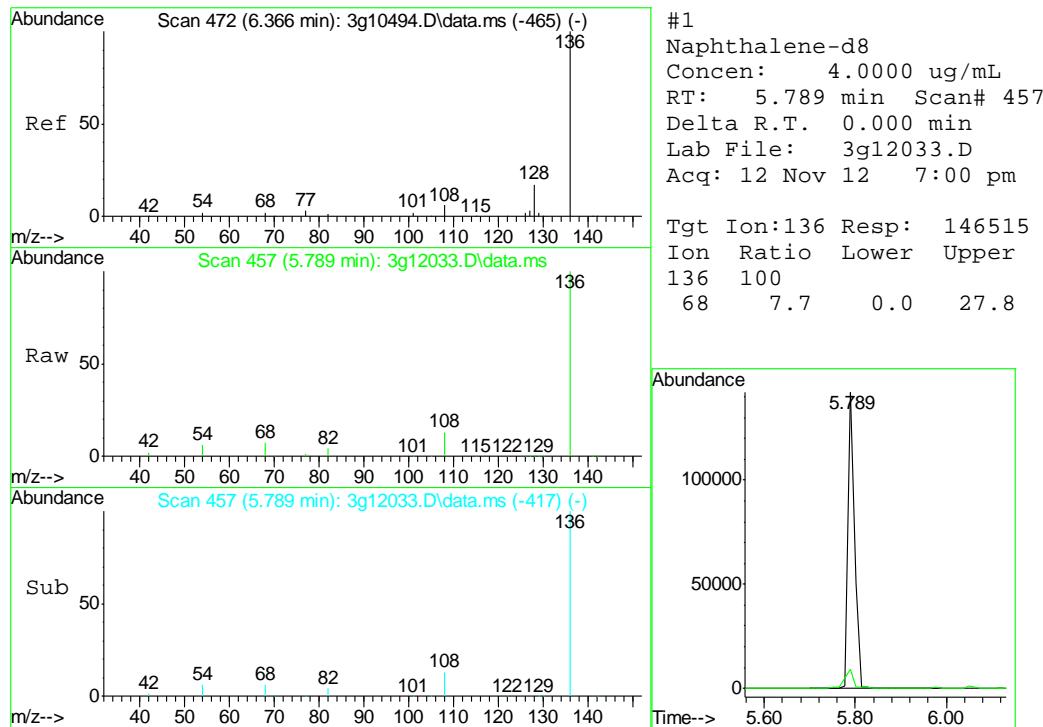
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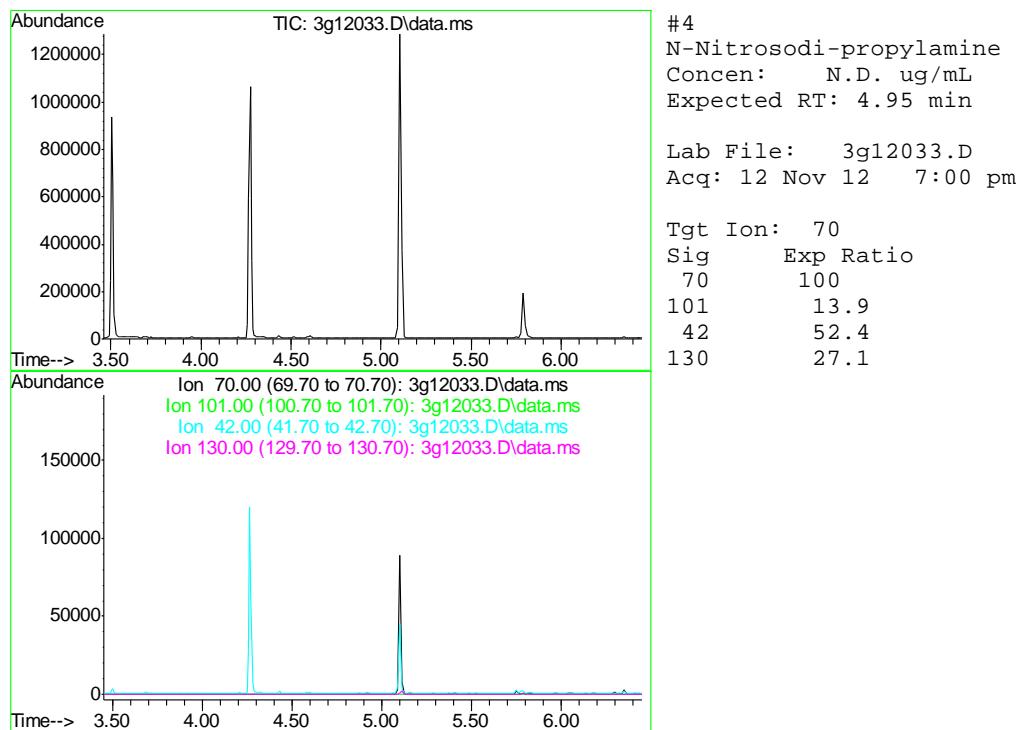
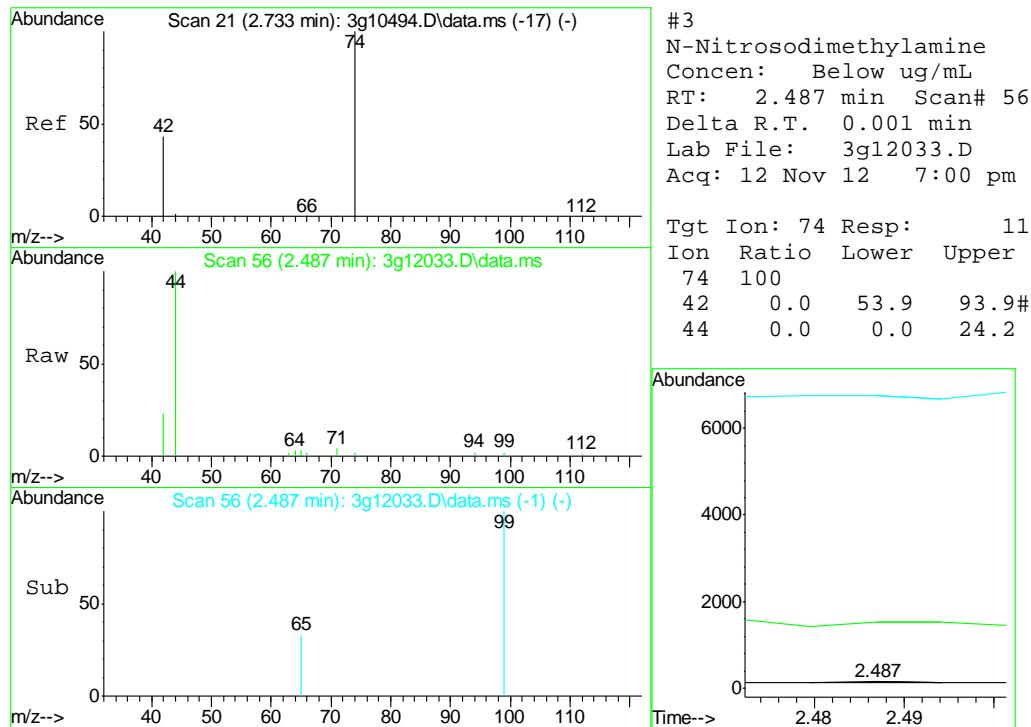
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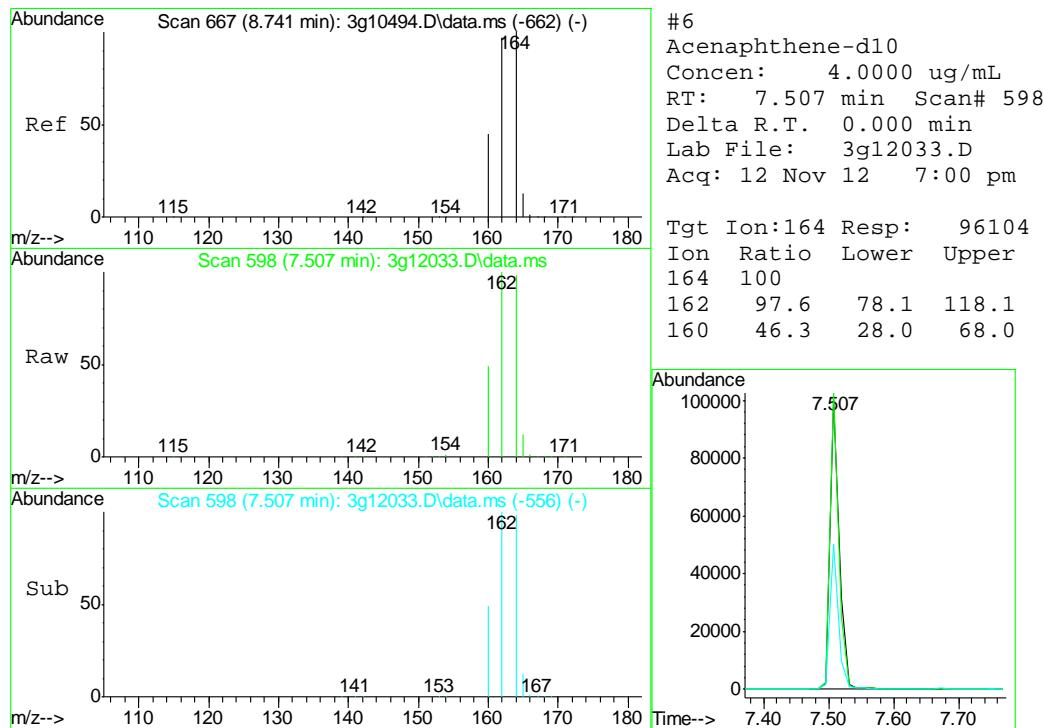
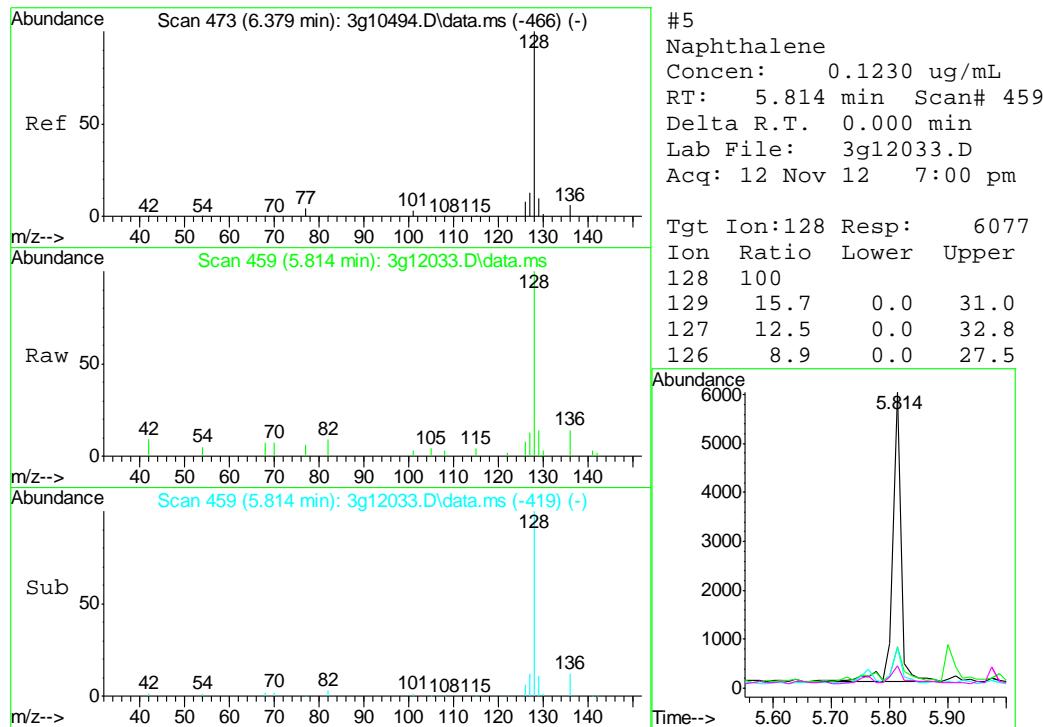
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 Operator : DONC  
 Sample : D40798-1  
 Misc : OP6947,E3G568,30.09,,,1,1  
 ALS Vial : 19 Sample Multiplier: 1

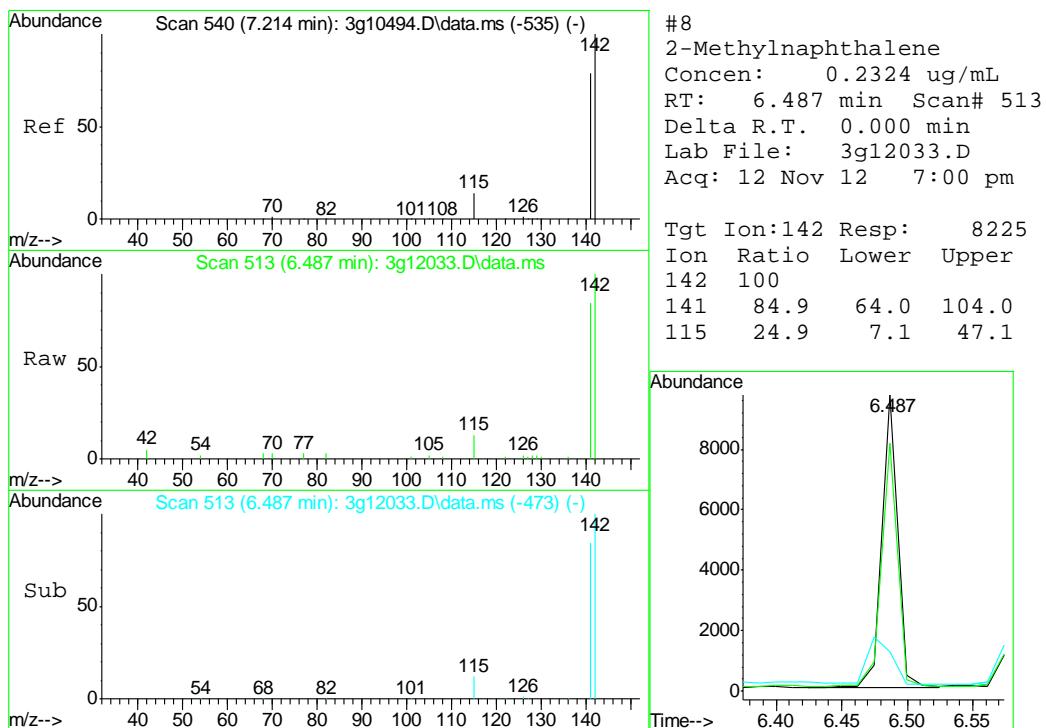
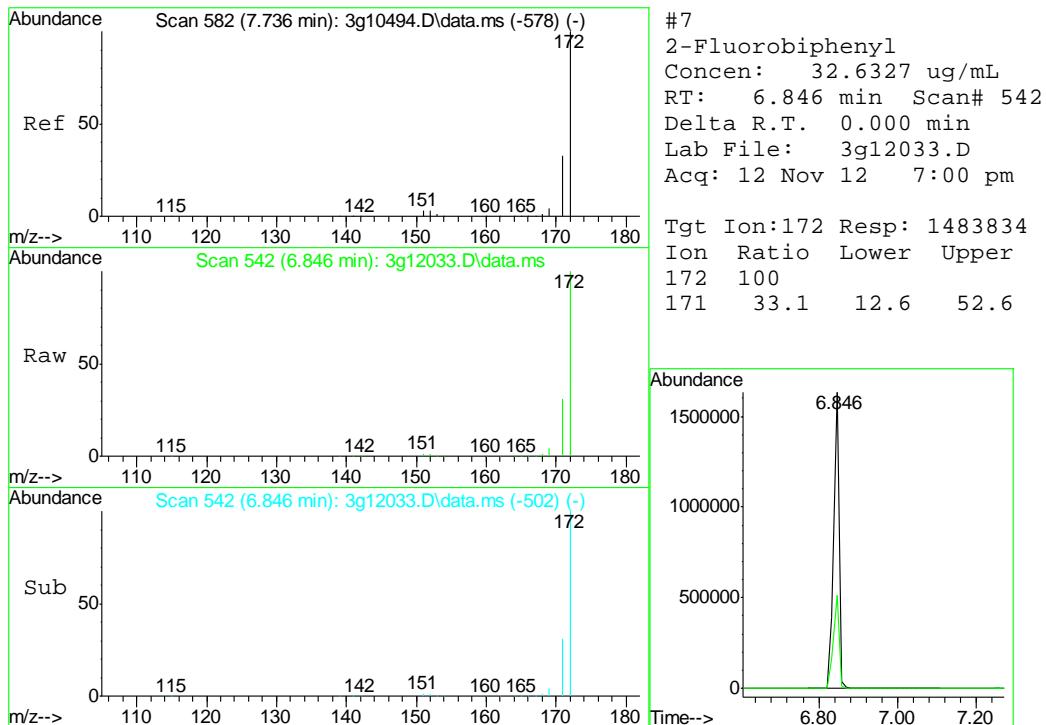
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 Quant Title : PAHSIM BASE  
 QLast Update : Mon Nov 12 15:44:05 2012  
 Response via : Initial Calibration

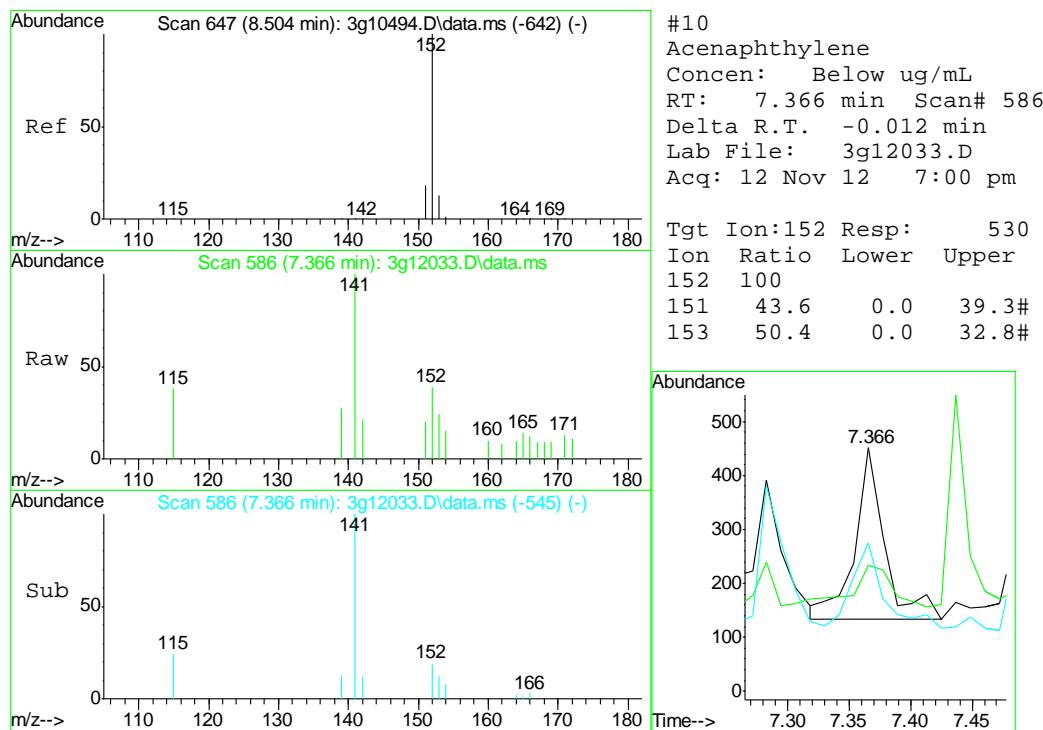
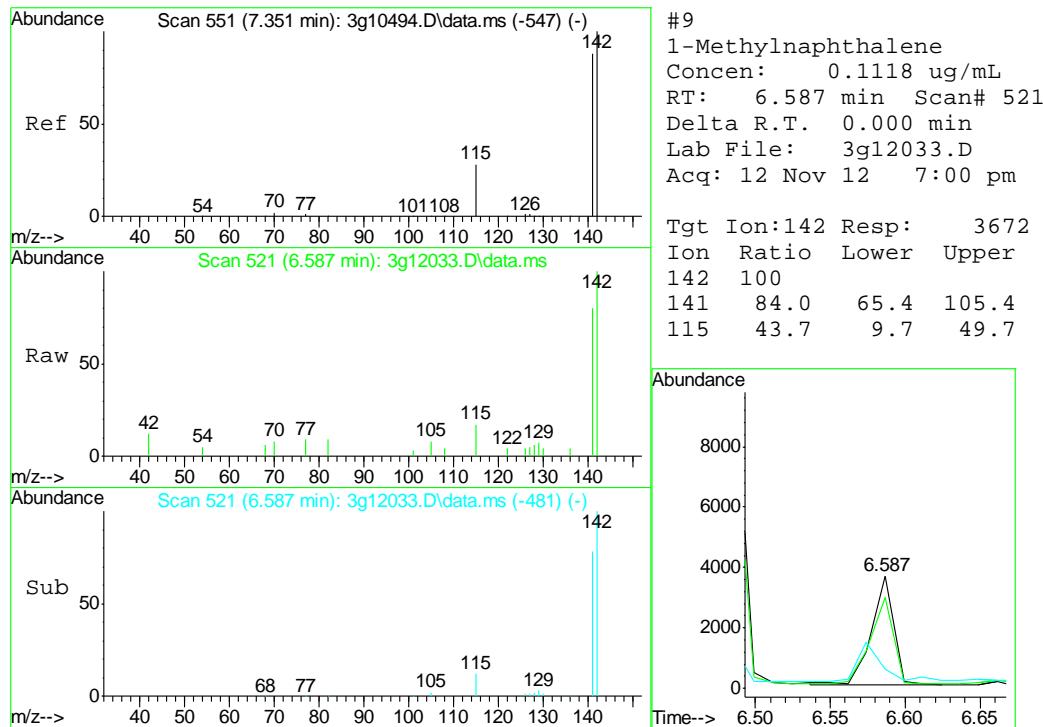


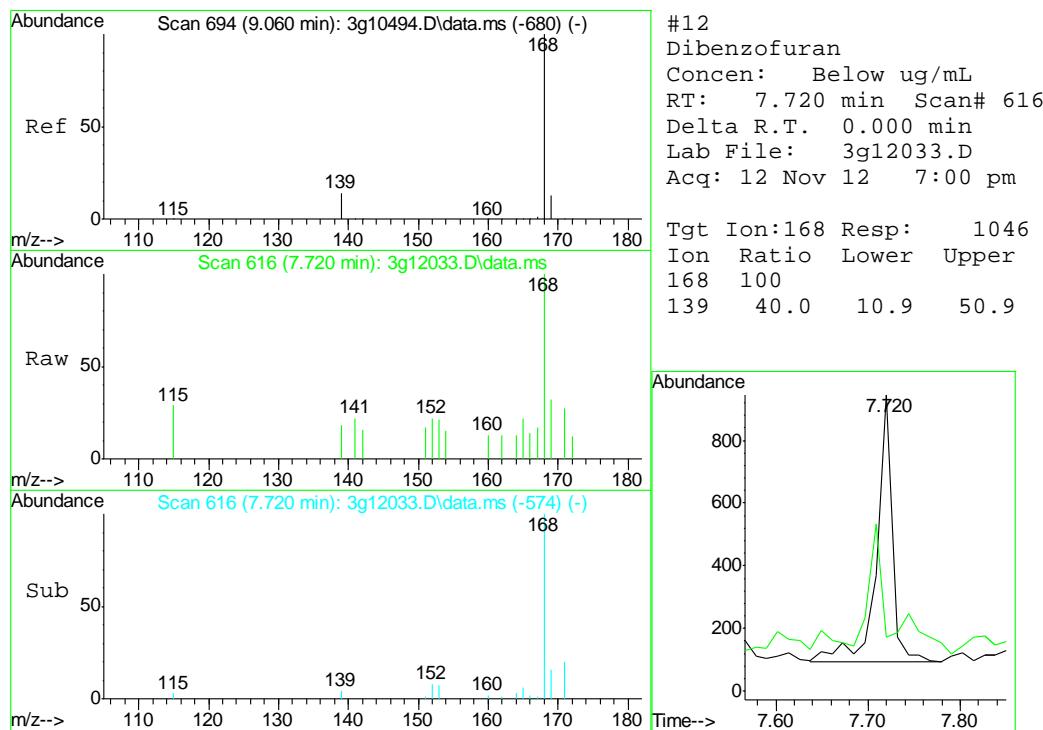
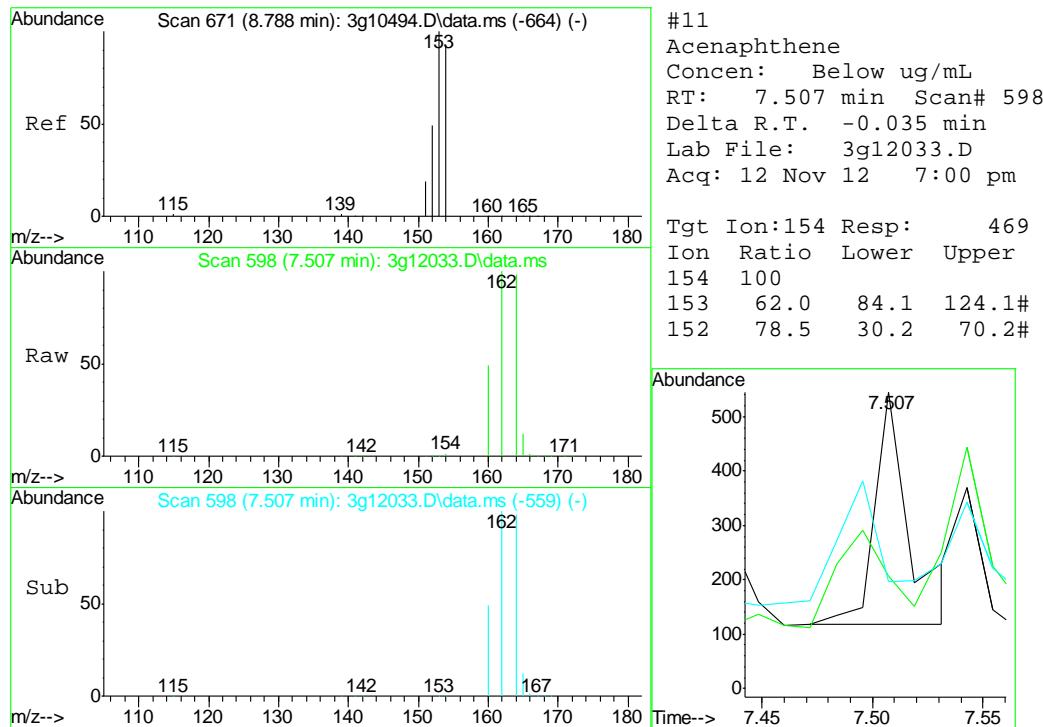


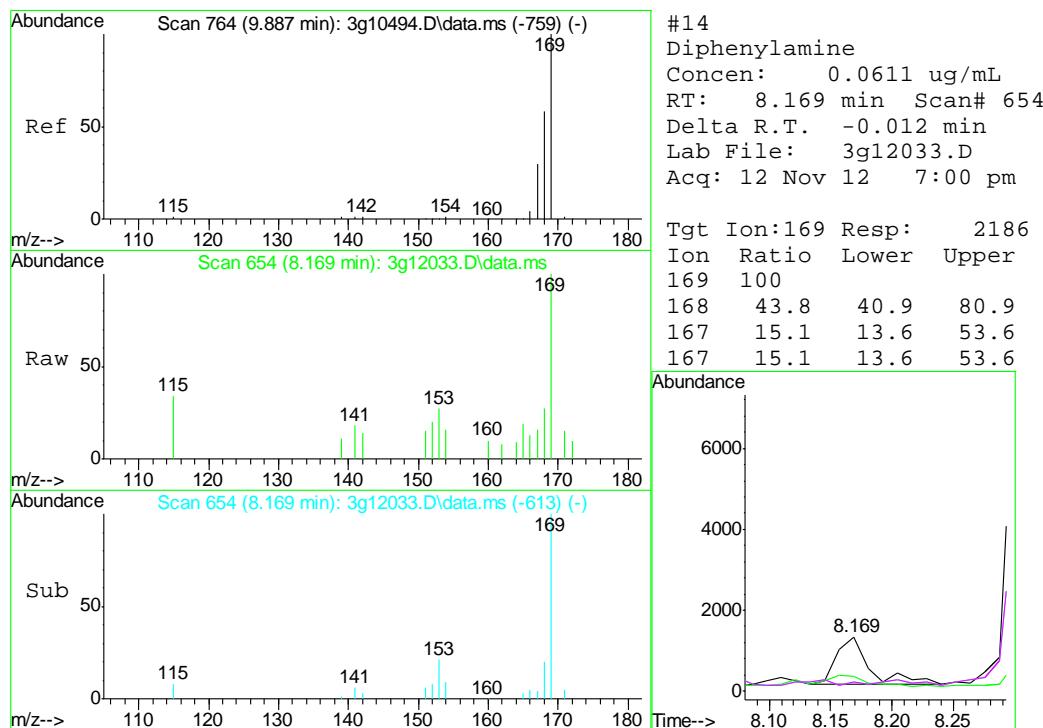
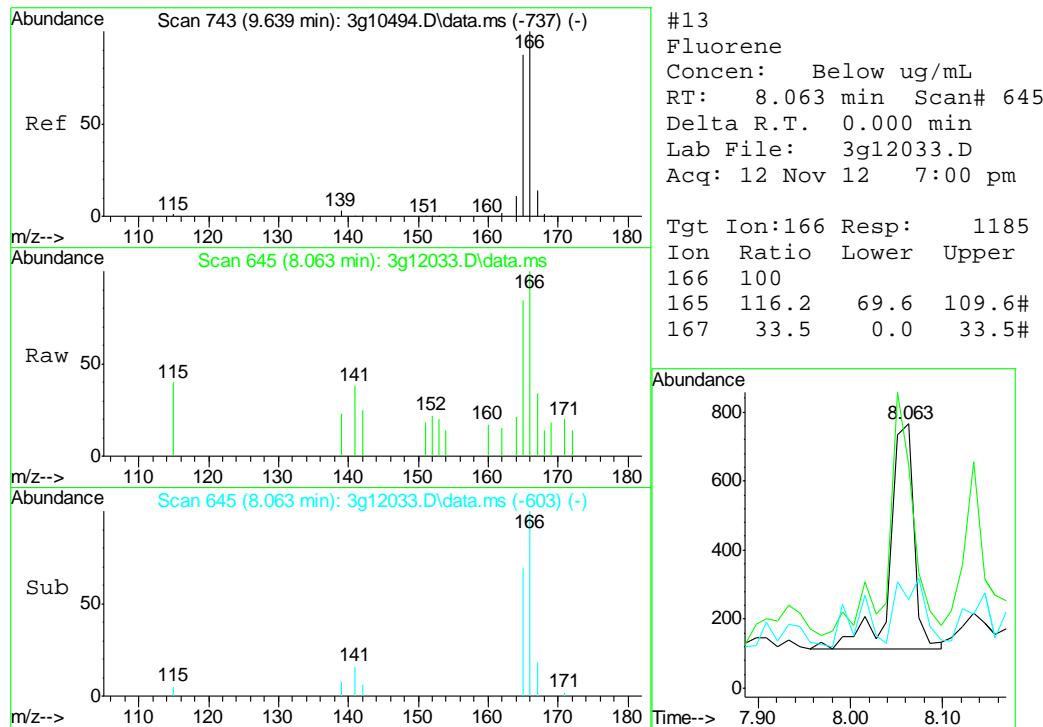


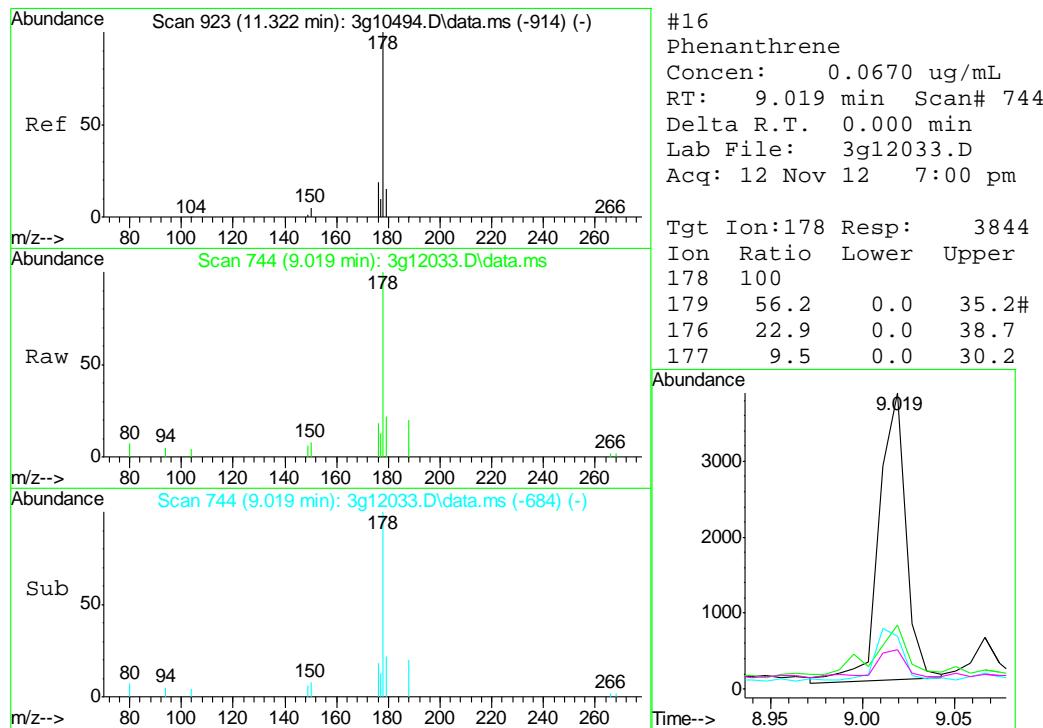
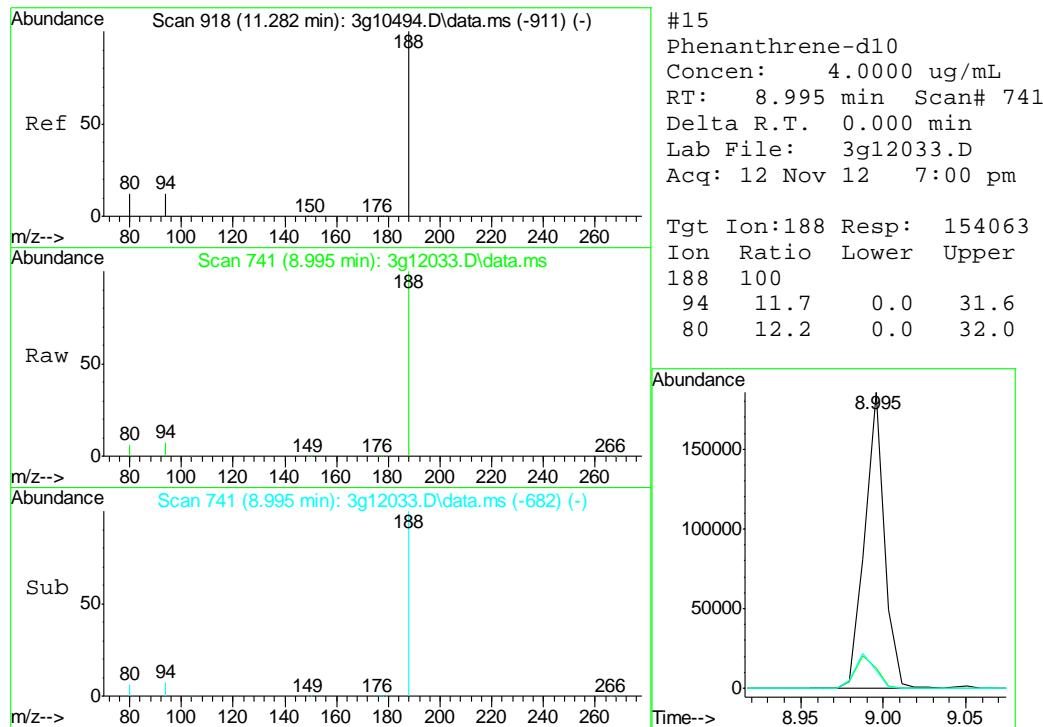


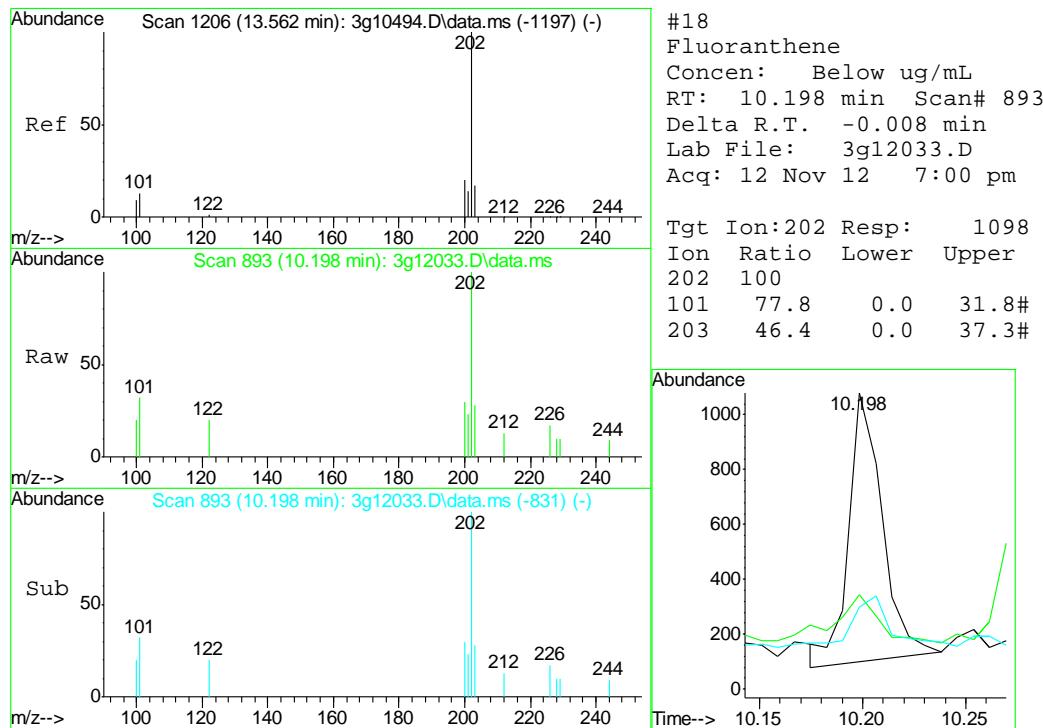
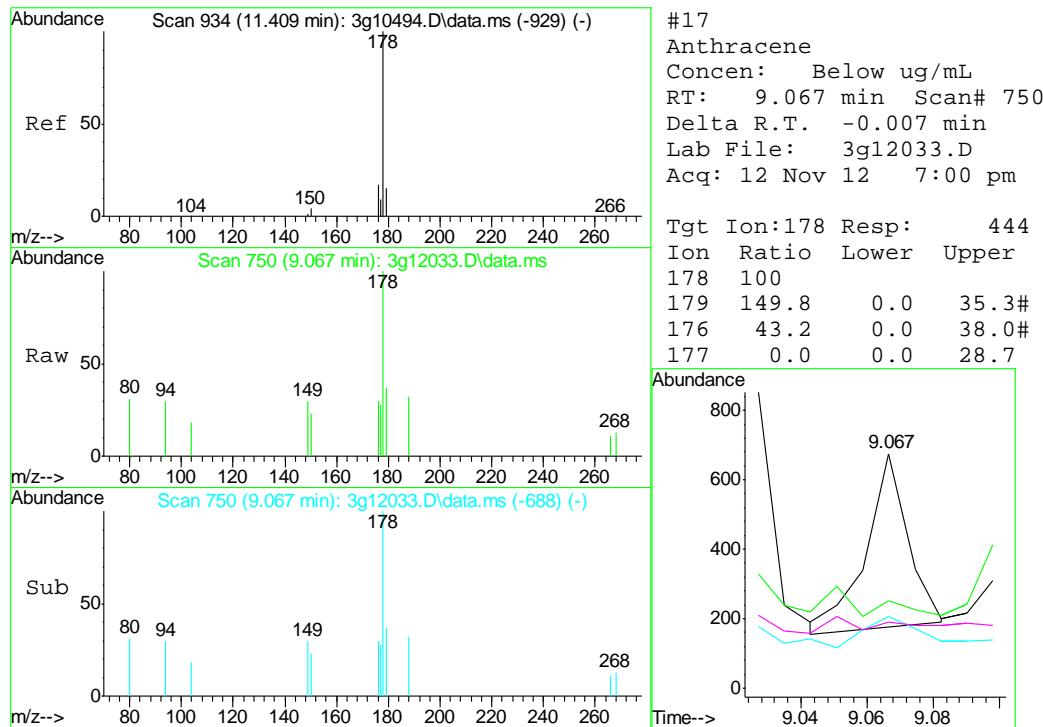


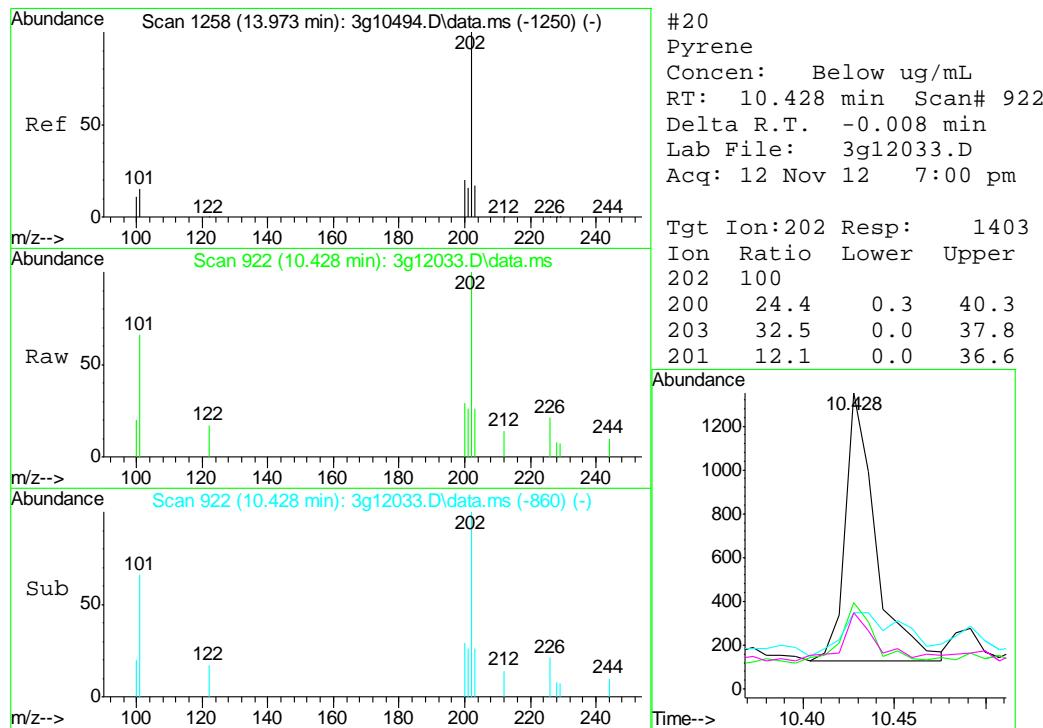
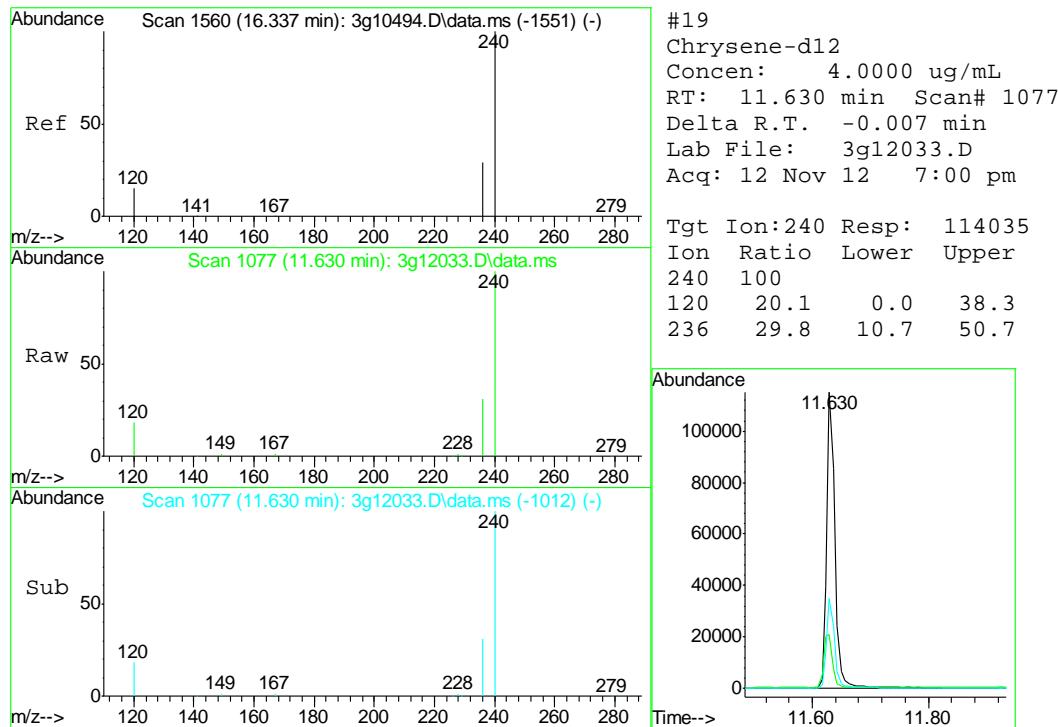


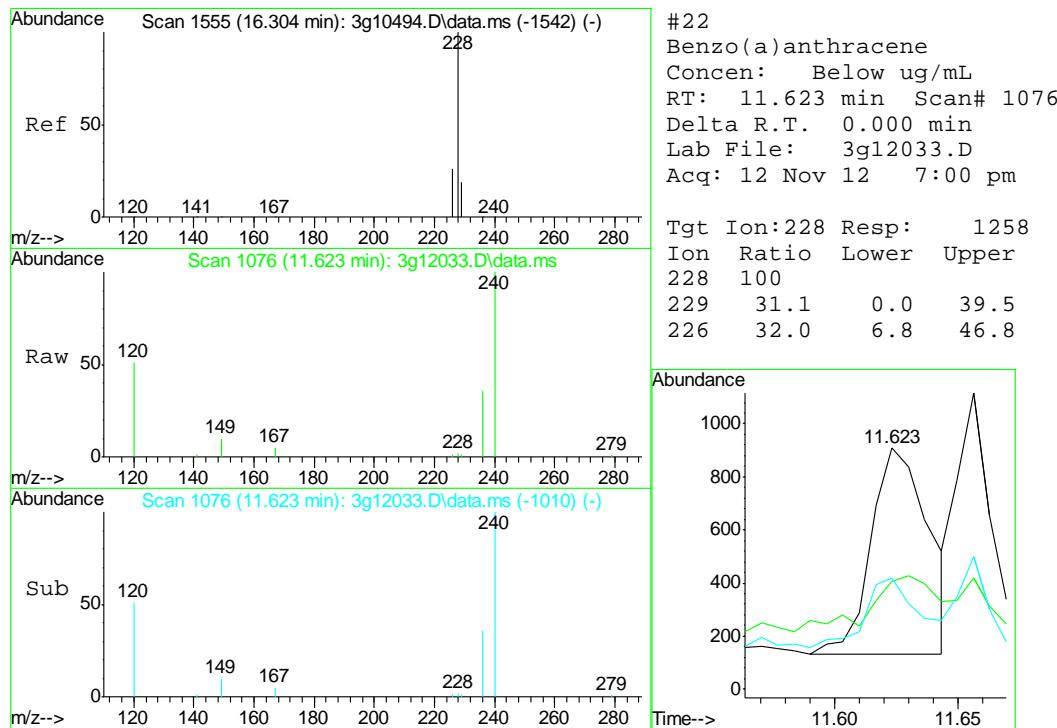
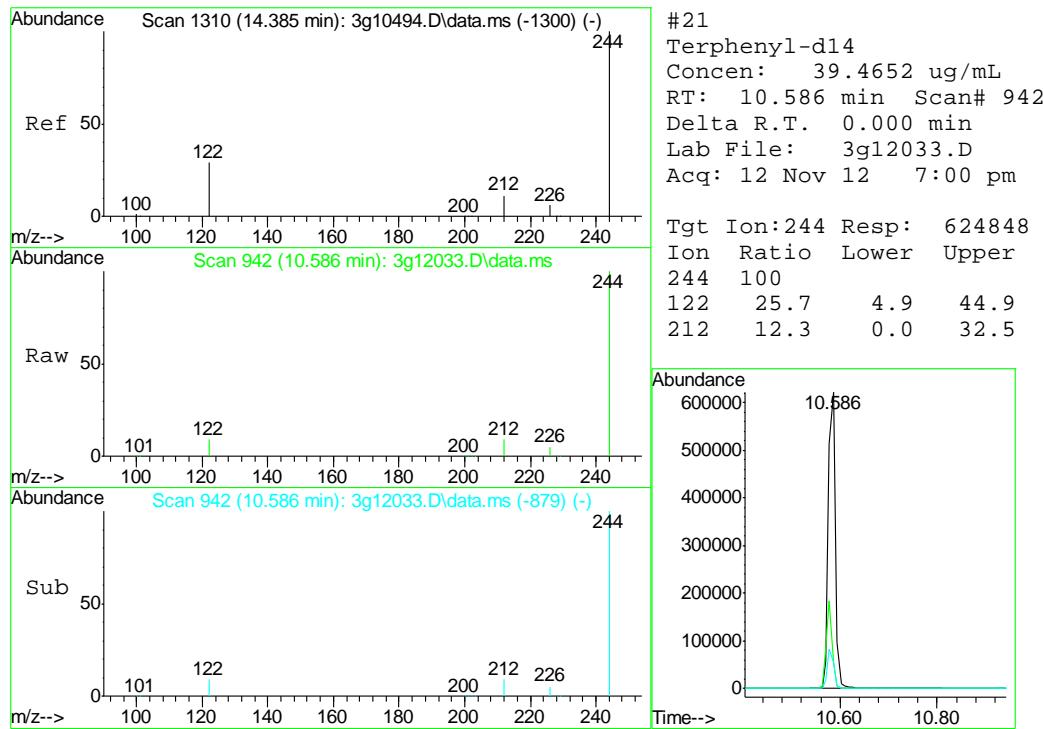


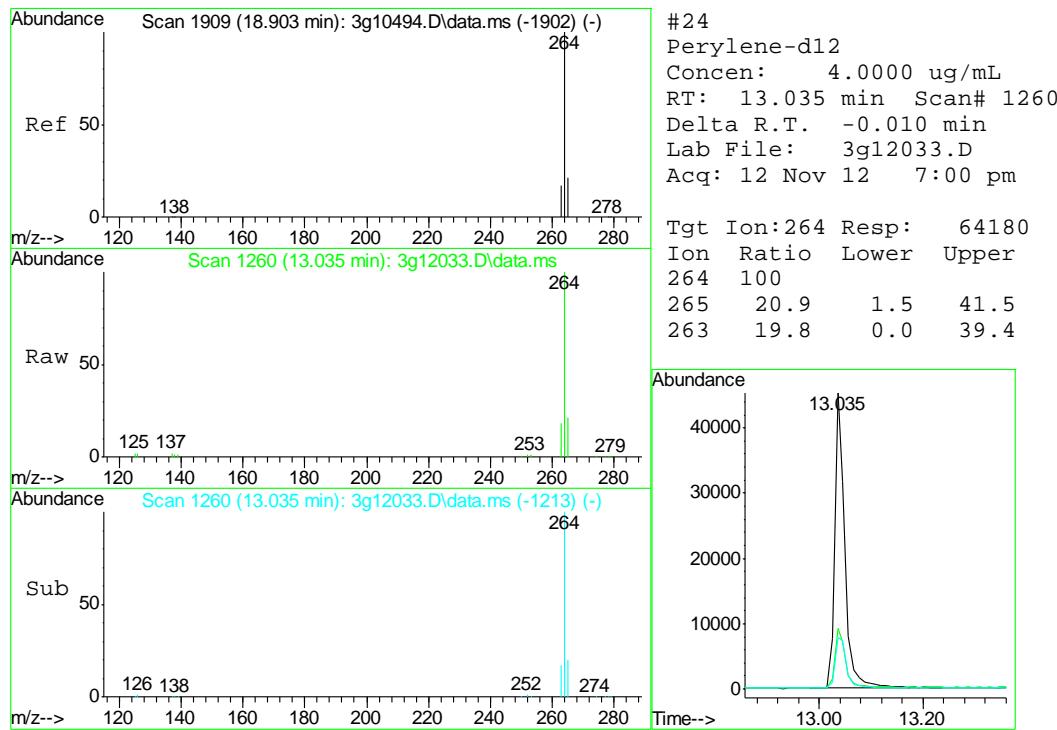
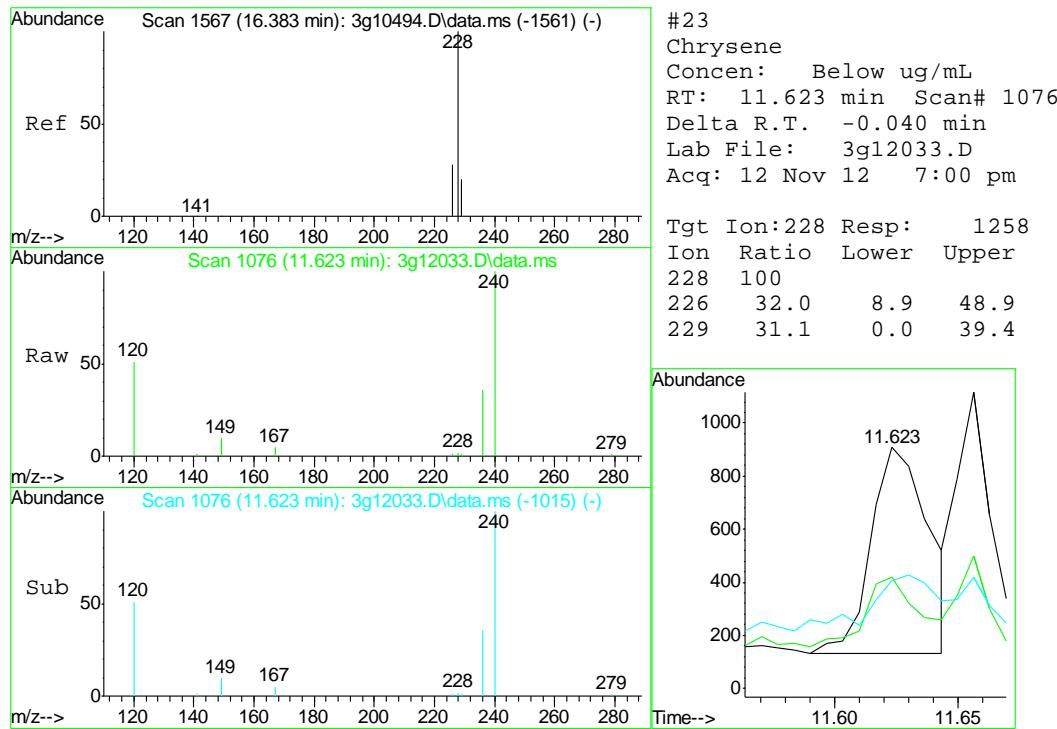


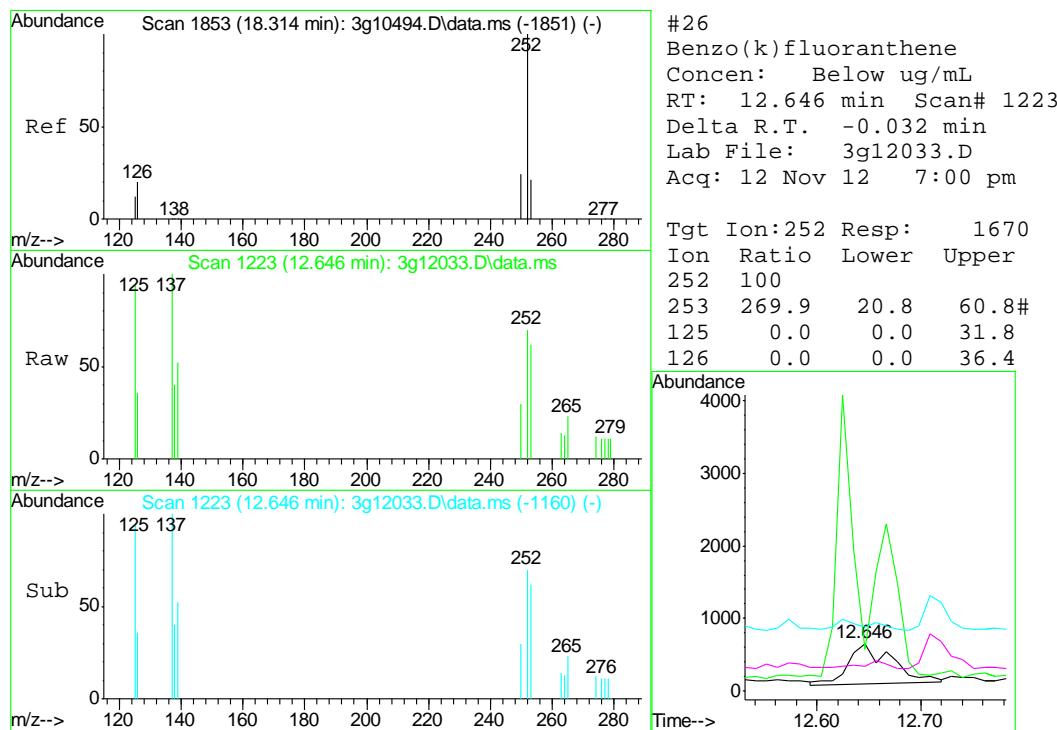
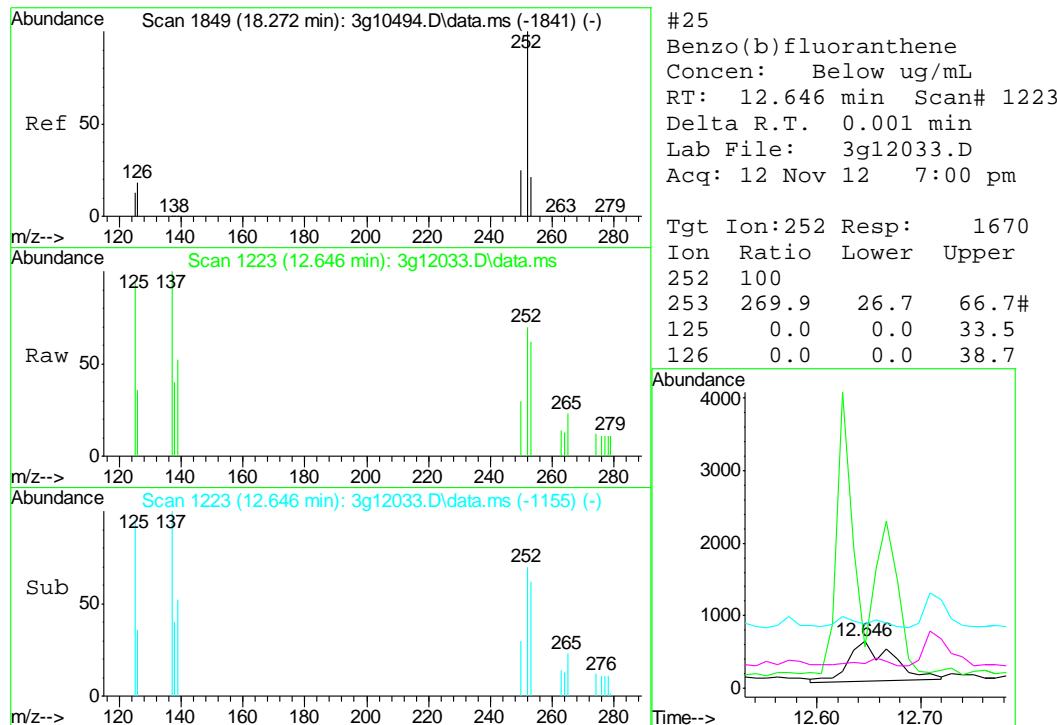


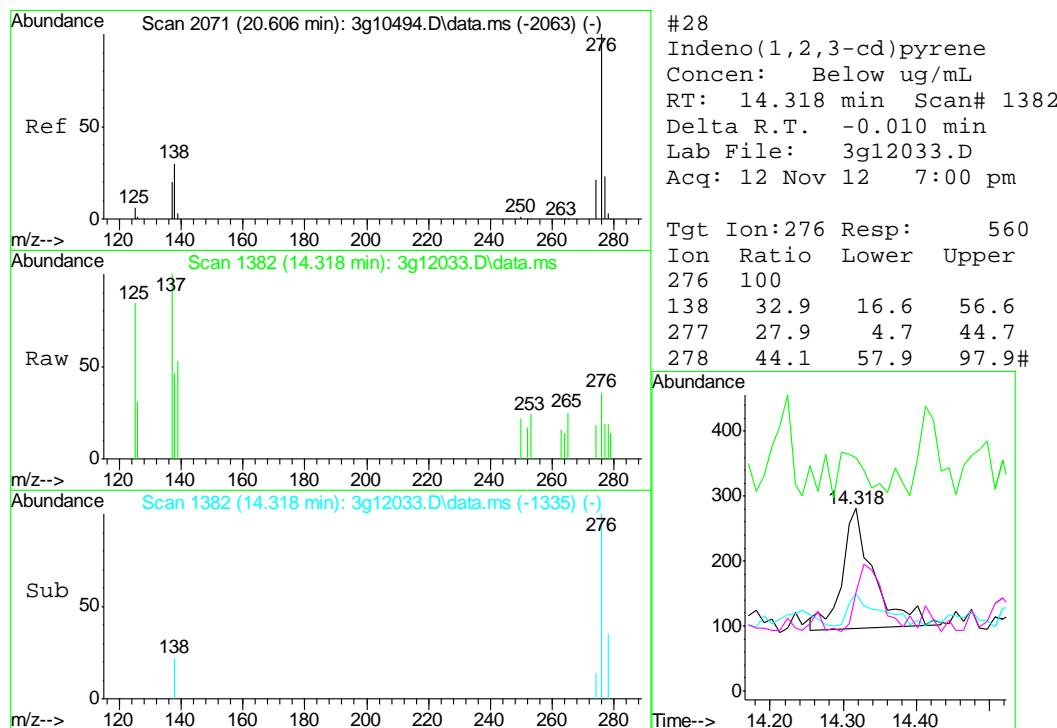
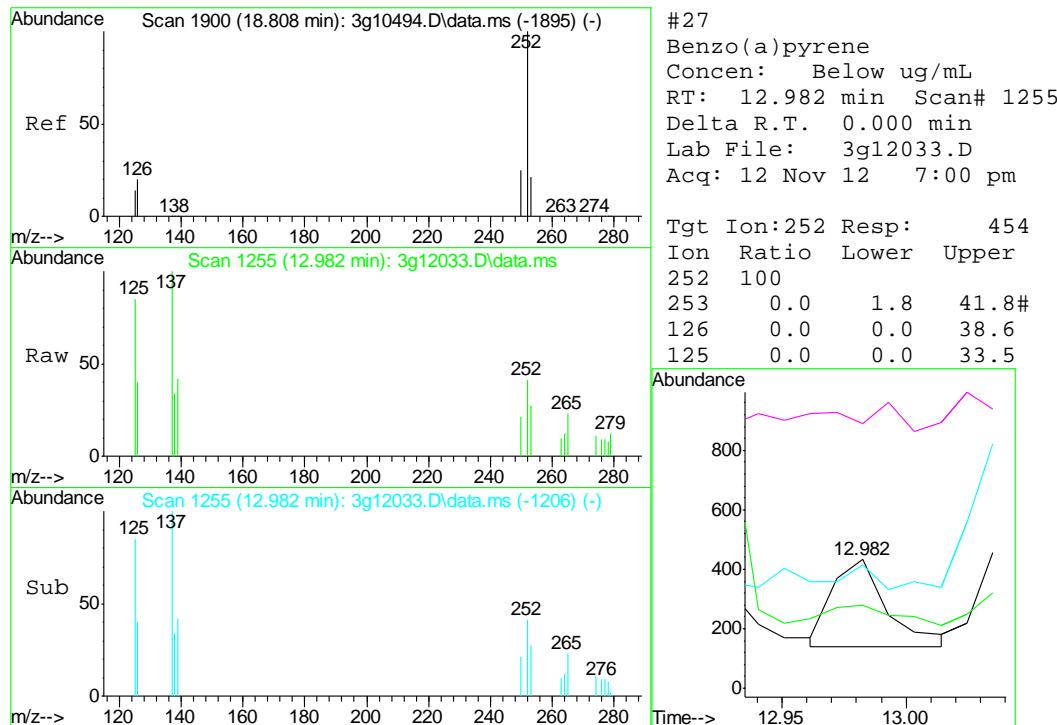


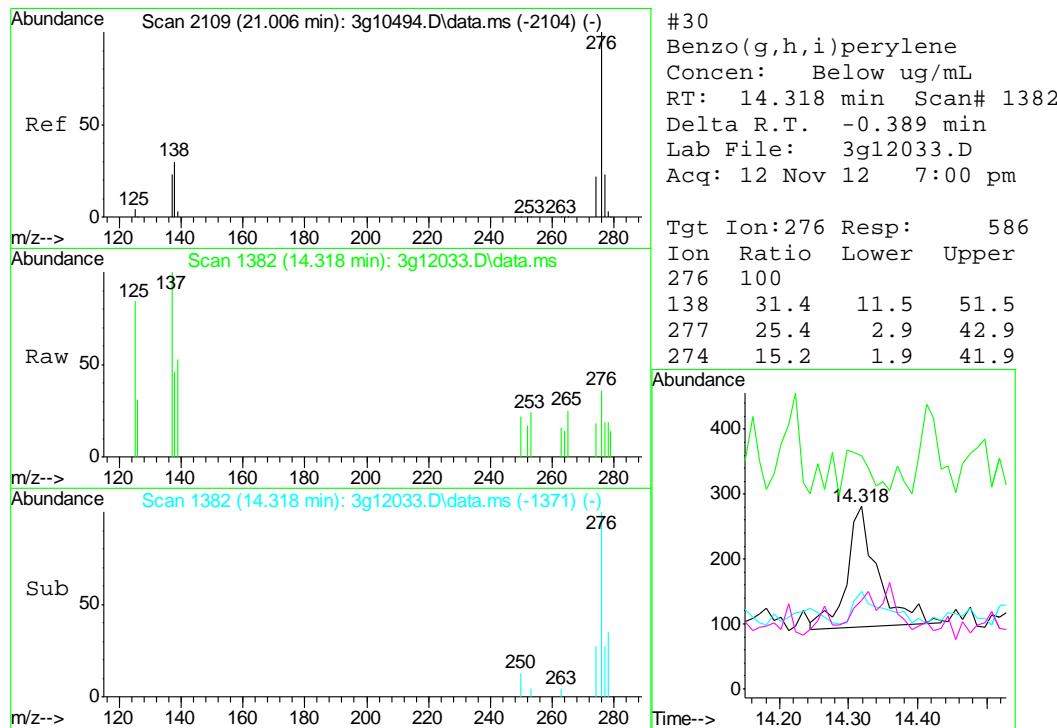
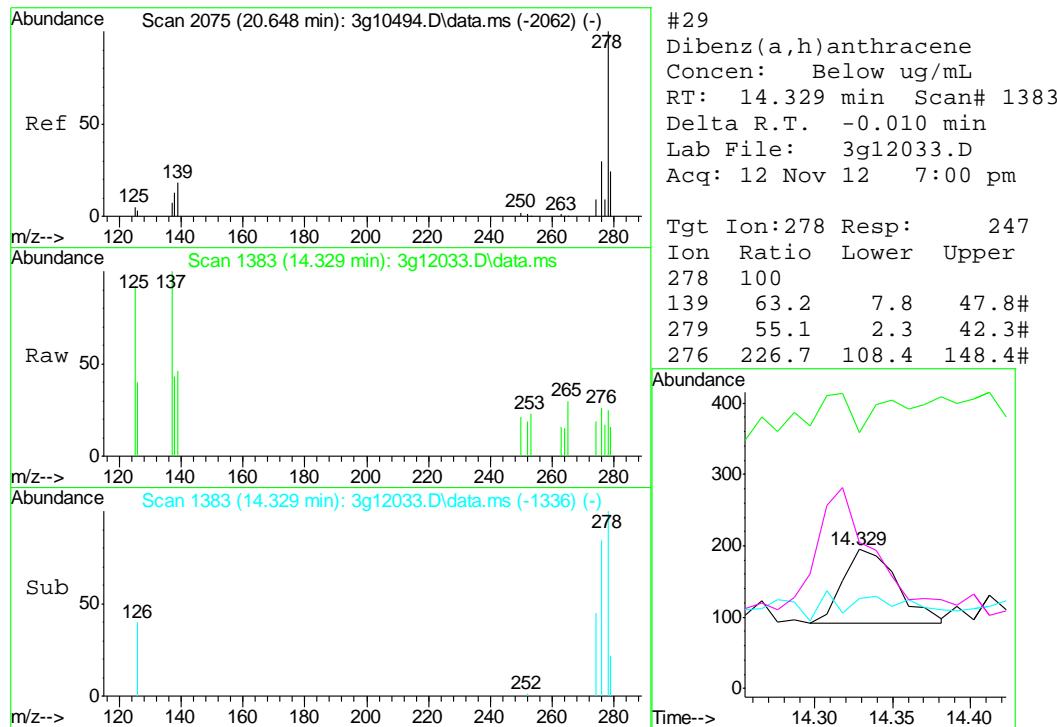












## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\111212\  
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 Operator : DONC  
 Sample : OP6947-MB  
 Misc : OP6947,E3G568,30.00,,,1,1  
 ALS Vial : 12 Sample Multiplier: 1

Quant Time: Nov 12 16:40:27 2012  
 Quant Method : C:\msdchem\1\METHODS\SIMPE3G568.M  
 Quant Title : PAHSIM BASE  
 QLast Update : Mon Nov 12 15:44:05 2012  
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Naphthalene-d8	5.789	136	144918	4.0000	ug/mL	0.00
6) Acenaphthene-d10	7.507	164	90962	4.0000	ug/mL	0.00
15) Phenanthrene-d10	8.995	188	157997	4.0000	ug/mL	0.00
19) Chrysene-d12	11.636	240	113918	4.0000	ug/mL	0.00
24) Perylene-d12	13.046	264	63583	4.0000	ug/mL	0.00

System Monitoring Compounds						
2) Nitrobenzene-d5	5.103	82	608123	42.1273	ug/mL	0.00
Spiked Amount	50.000	Range	25 - 135	Recovery	= 84.26%	
7) 2-Fluorobiphenyl	6.846	172	1730035	40.1979	ug/mL	0.00
Spiked Amount	50.000	Range	25 - 135	Recovery	= 80.40%	
21) Terphenyl-d14	10.586	244	741133	46.8578	ug/mL	0.00
Spiked Amount	50.000	Range	25 - 135	Recovery	= 93.72%	

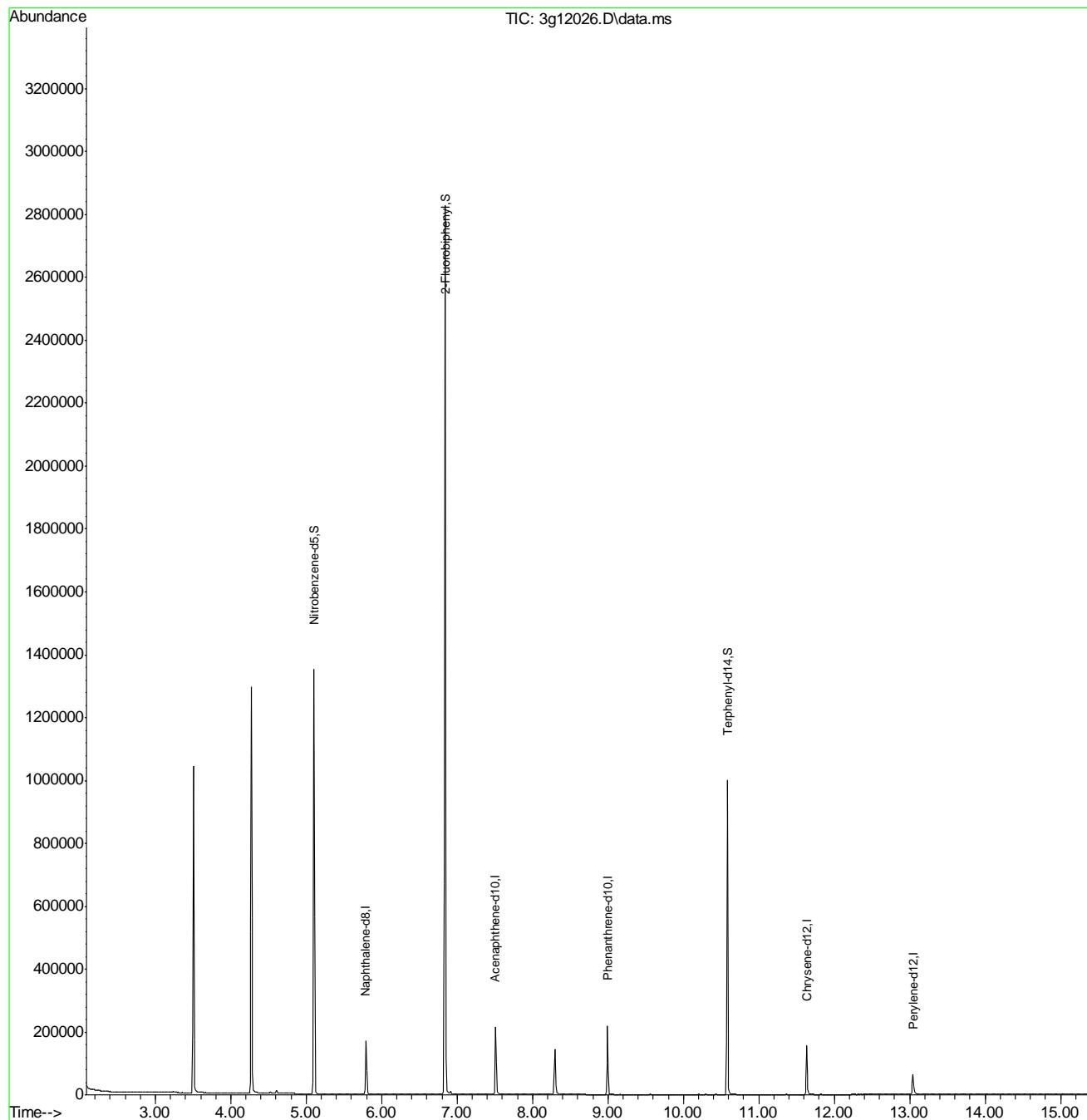
Target Compounds					Qvalue
3) N-Nitrosodimethylamine	2.501	74	25	N.D.	
4) N-Nitrosodi-propylamine	0.000	70	0	N.D. d	
5) Naphthalene	5.814	128	847	N.D.	
8) 2-Methylnaphthalene	6.487	142	836	N.D.	
9) 1-Methylnaphthalene	6.487	142	807	N.D.	
10) Acenaphthylene	7.377	152	1014	N.D.	
11) Acenaphthene	7.507	154	537	Below Cal #	89
12) Dibenzofuran	7.720	168	739	N.D.	
13) Fluorene	8.063	166	526	N.D.	
14) Diphenylamine	0.000	169	0	N.D. d	
16) Phenanthrene	9.067	178	476	N.D.	
17) Anthracene	9.067	178	476	N.D.	
18) Fluoranthene	10.206	202	568	N.D.	
20) Pyrene	10.436	202	566	N.D.	
22) Benzo(a)anthracene	11.630	228	1000	N.D.	
23) Chrysene	11.630	228	1000	N.D.	
25) Benzo(b)fluoranthene	12.646	252	794	N.D.	
26) Benzo(k)fluoranthene	12.646	252	794	N.D.	
27) Benzo(a)pyrene	12.646	252	794	N.D.	
28) Indeno(1,2,3-cd)pyrene	14.339	276	501	N.D.	
29) Dibenz(a,h)anthracene	14.339	278	252	N.D.	
30) Benzo(g,h,i)perylene	14.697	276	361	N.D.	

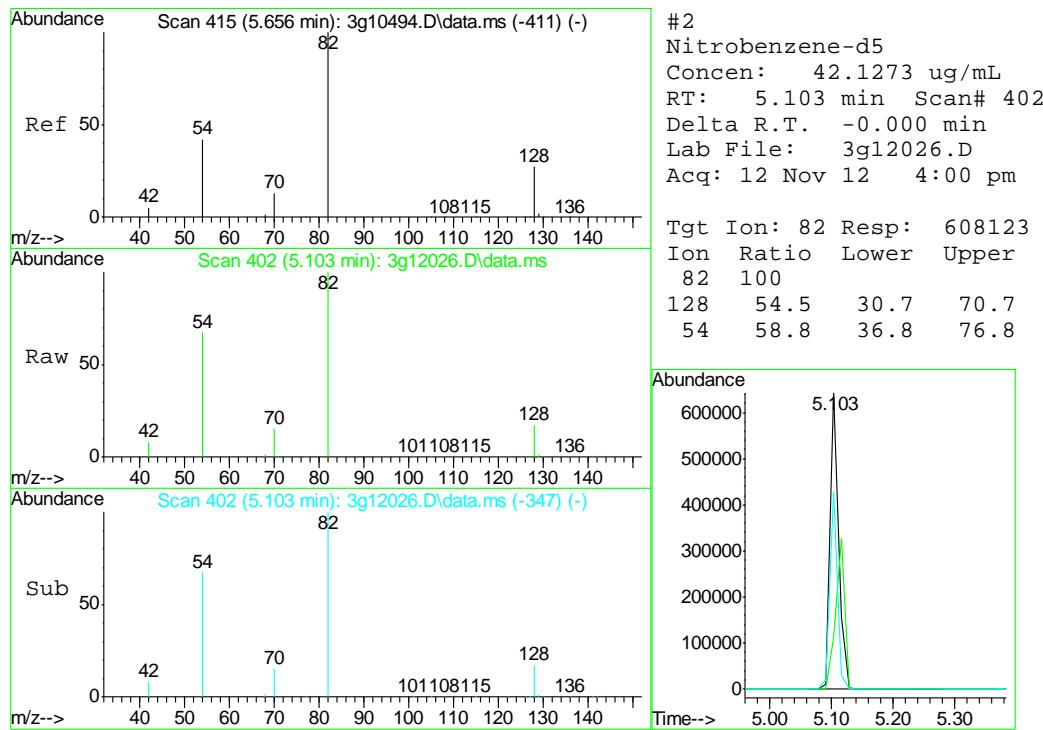
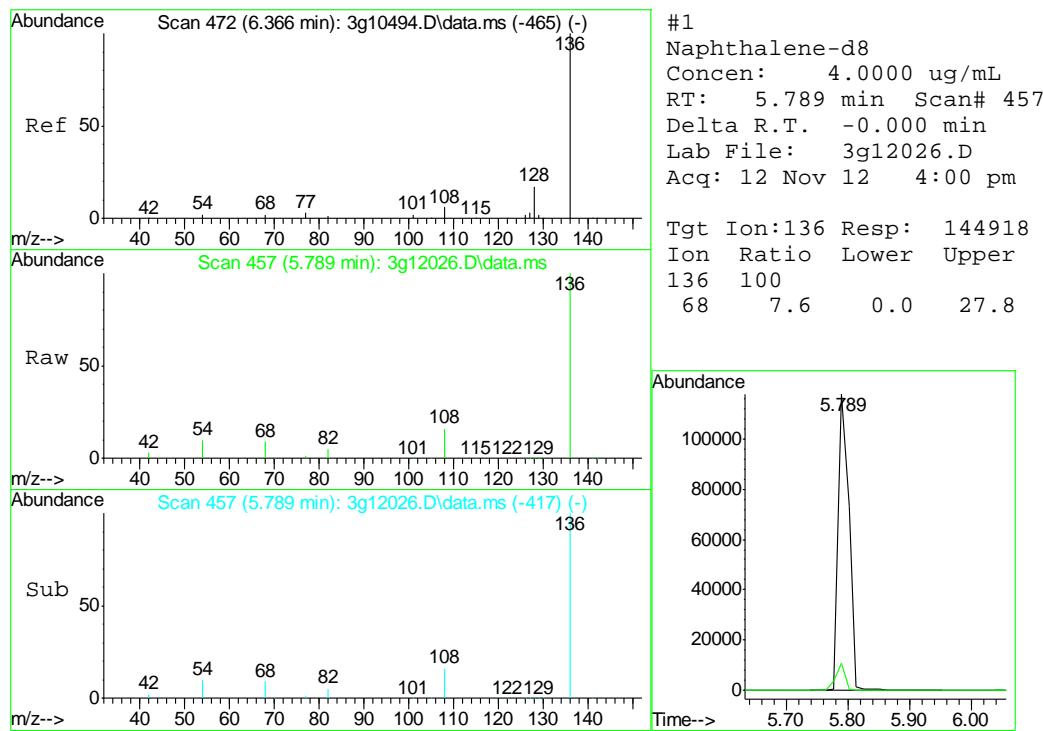
(#) = qualifier out of range (m) = manual integration (+) = signals summed

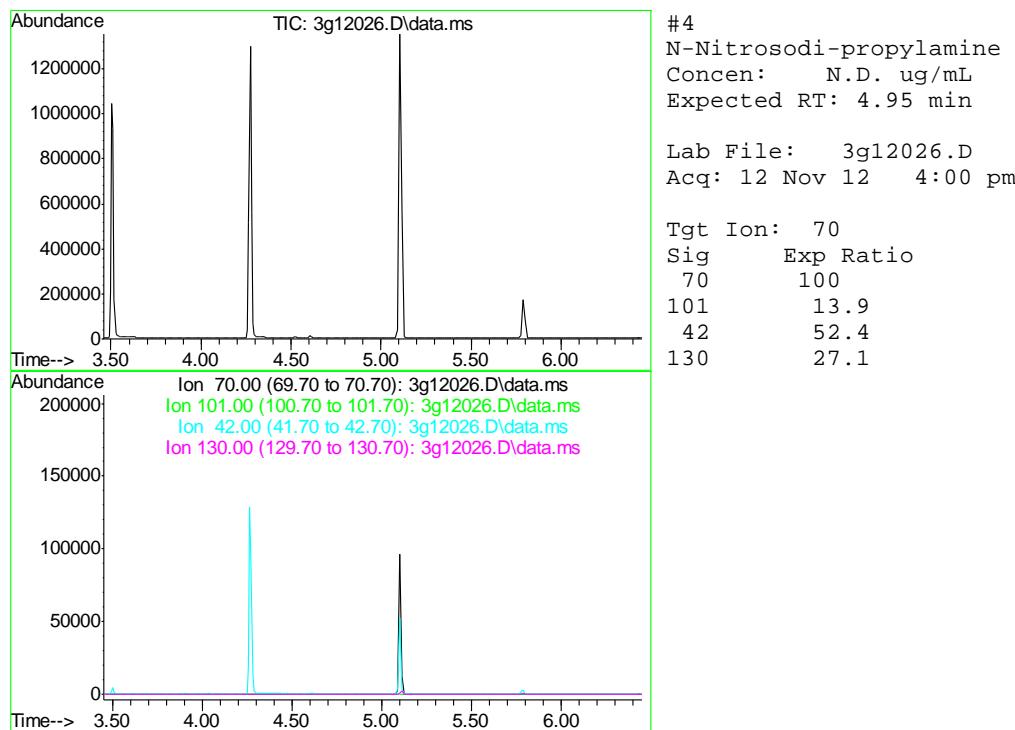
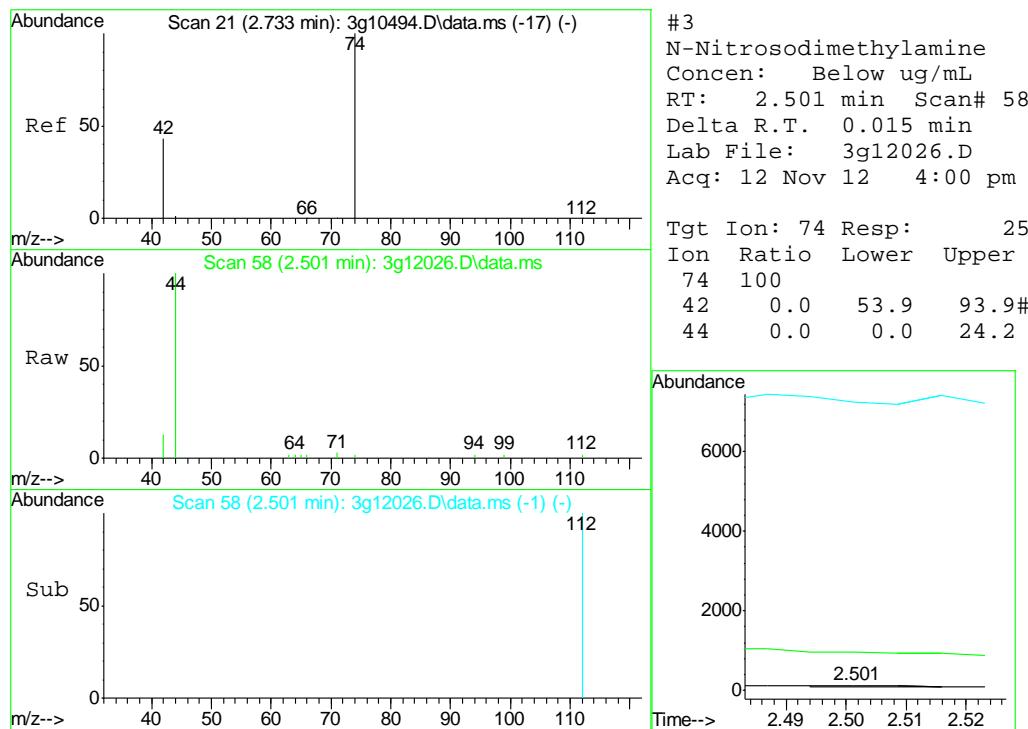
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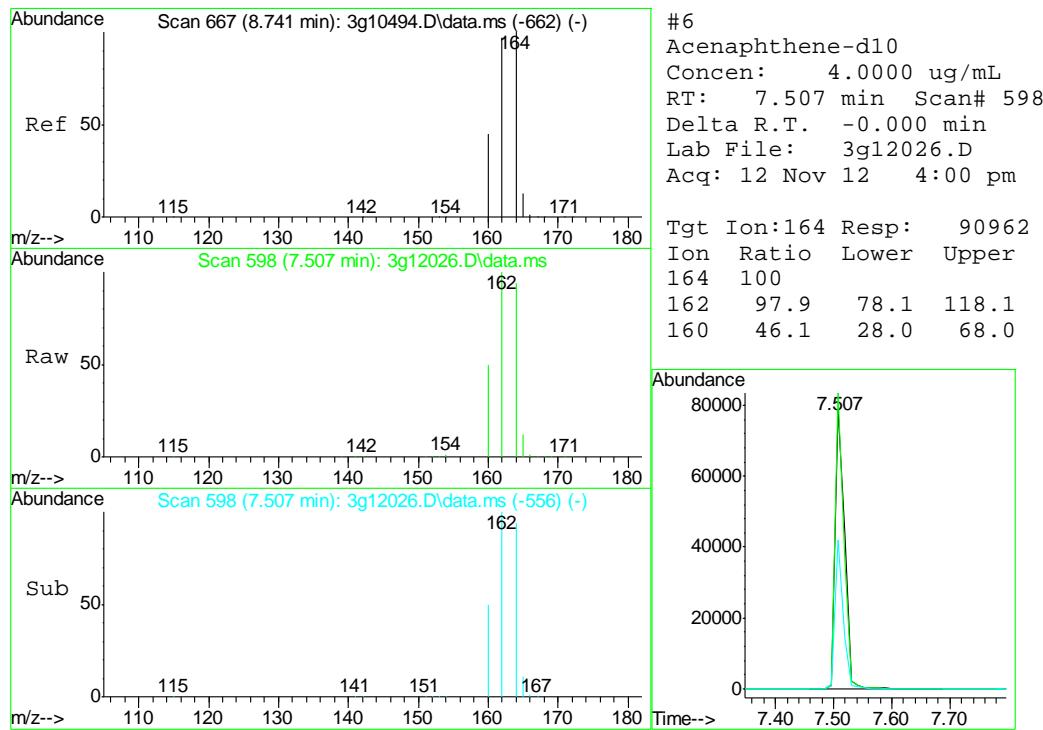
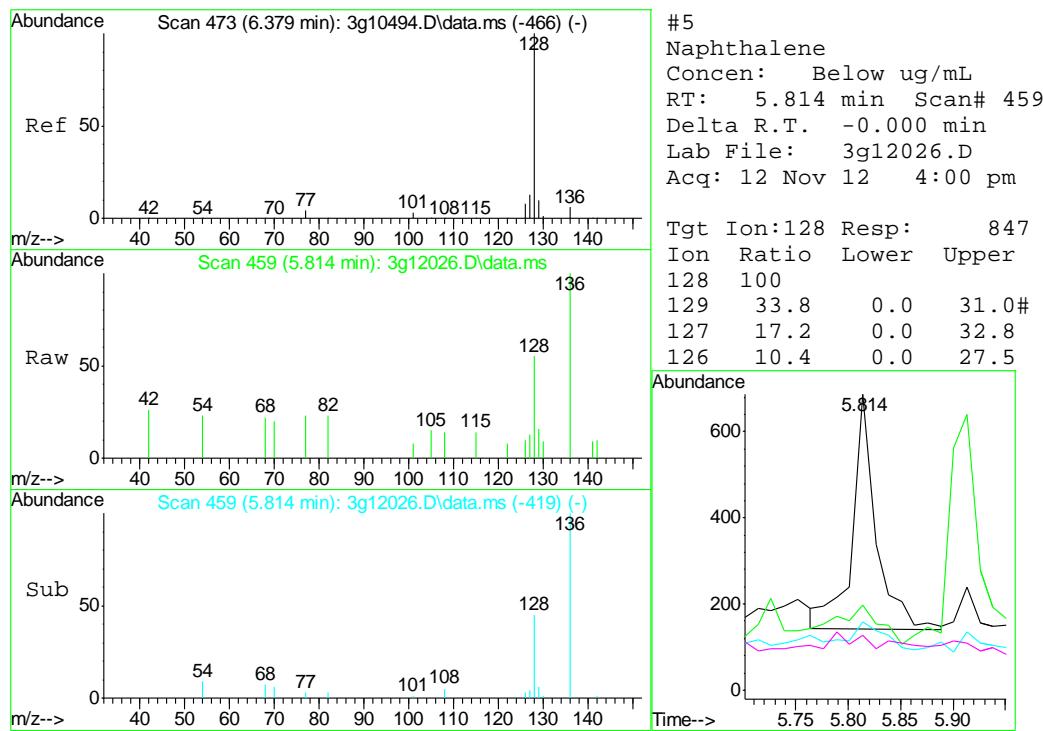
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 Sample : OP6947-MB  
 Misc : OP6947,E3G568,30.00,,,1,1  
 ALS Vial : 12 Sample Multiplier: 1

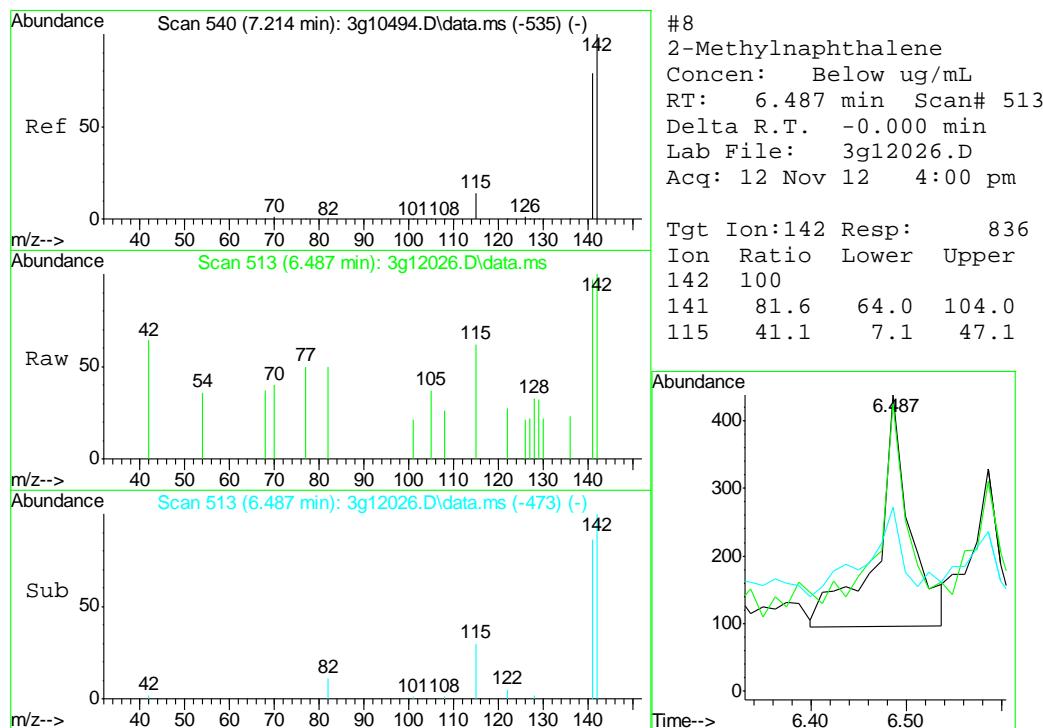
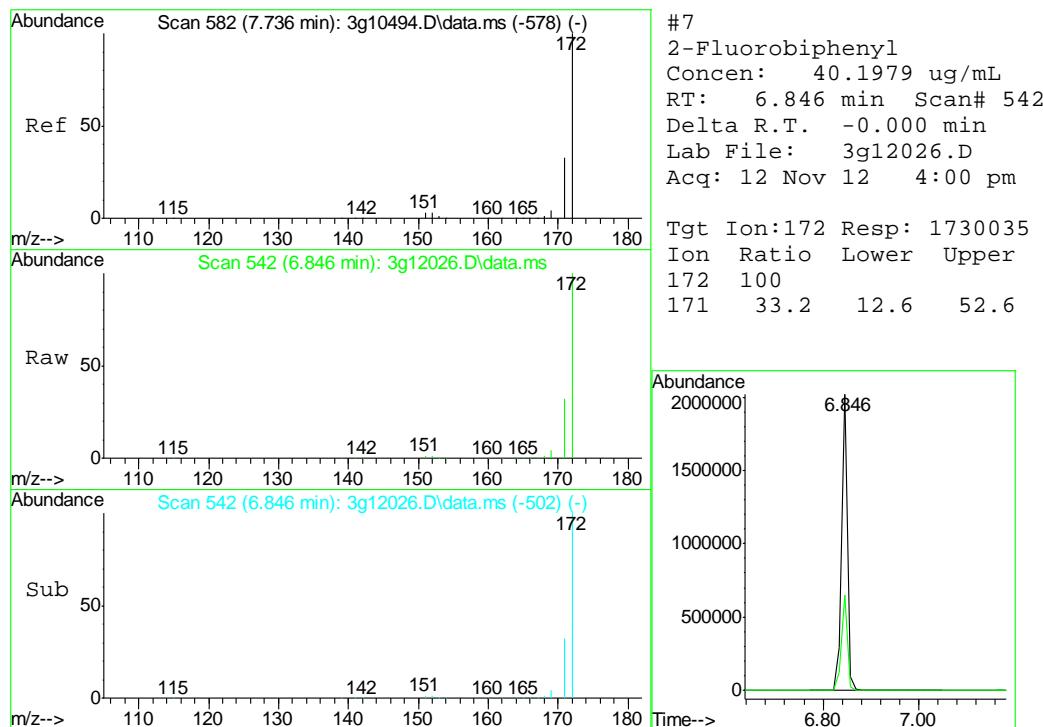
Quant Time: Nov 12 16:40:27 2012  
 Quant Method : C:\msdchem\1\METHODS\SIMPE3G568.M  
 Quant Title : PAHSIM BASE  
 QLast Update : Mon Nov 12 15:44:05 2012  
 Response via : Initial Calibration

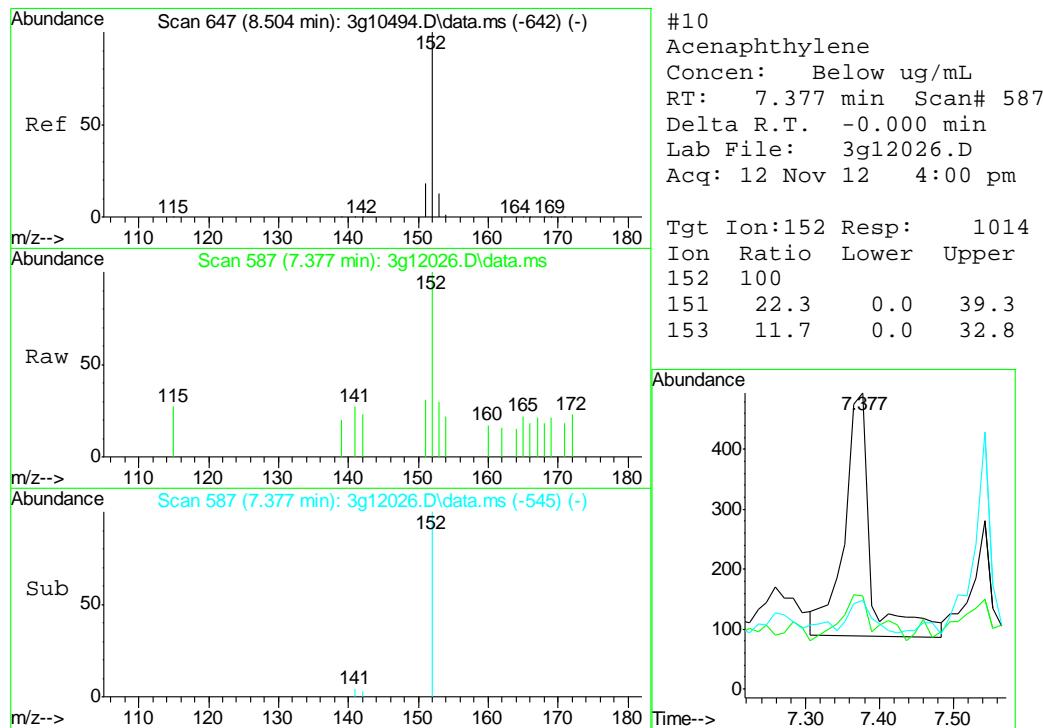
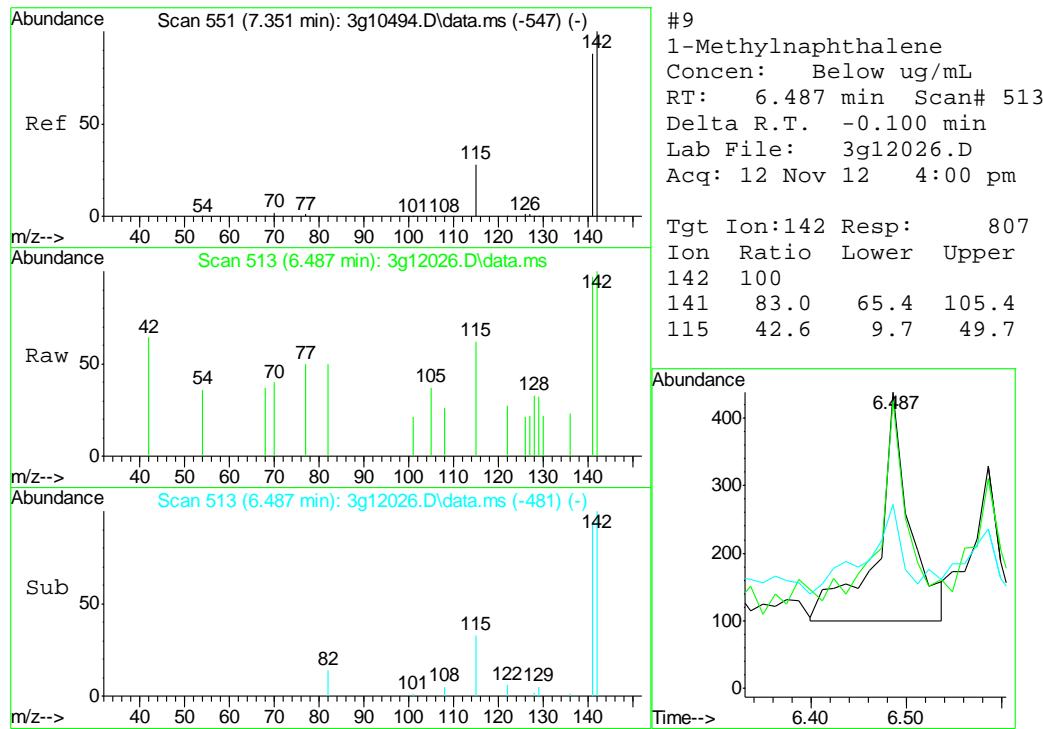


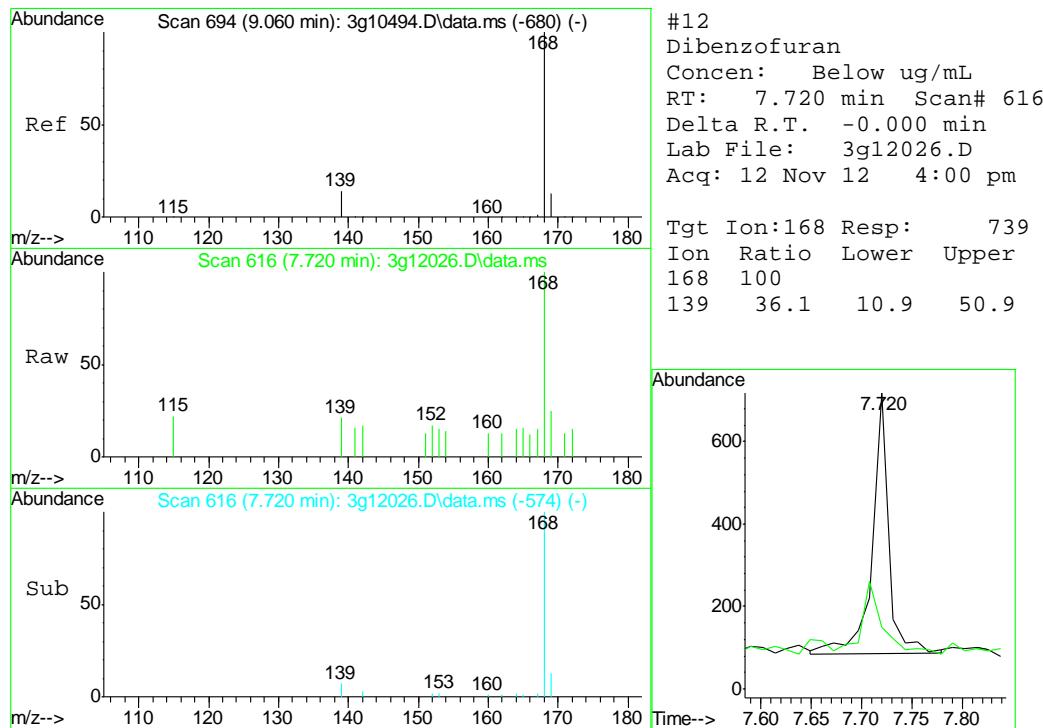
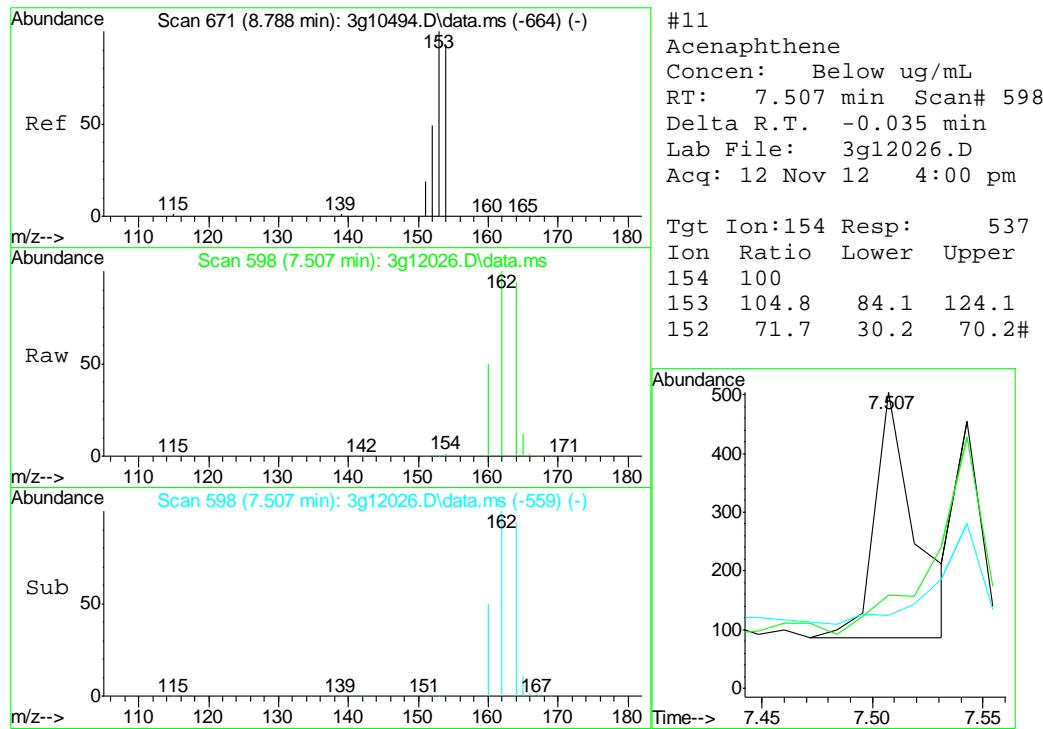


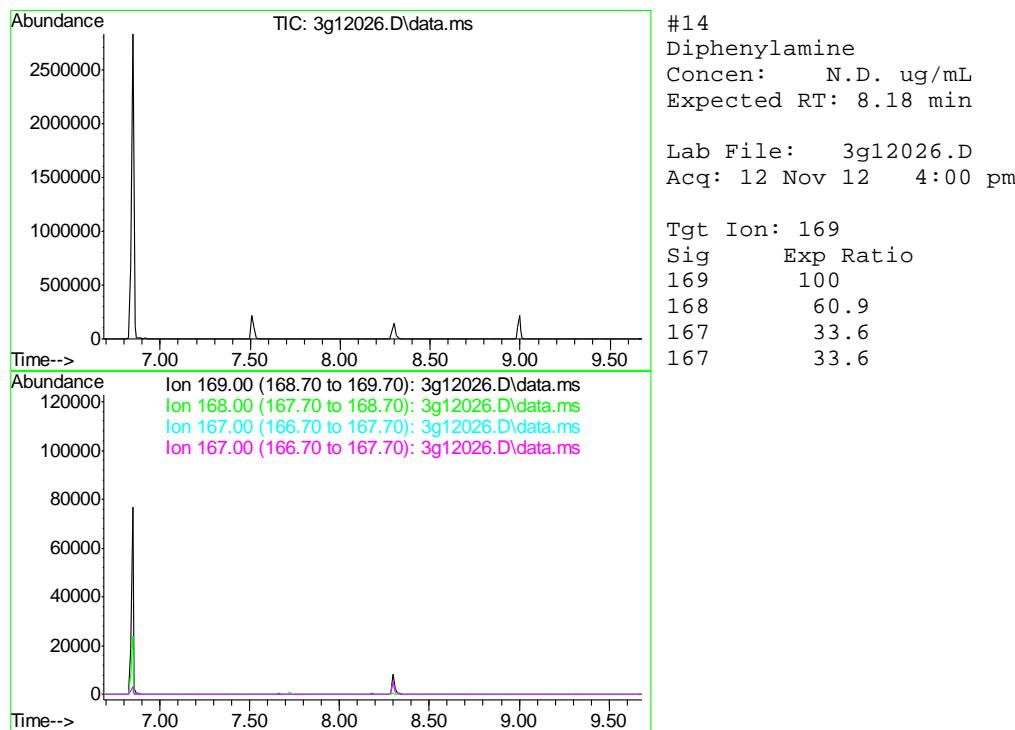
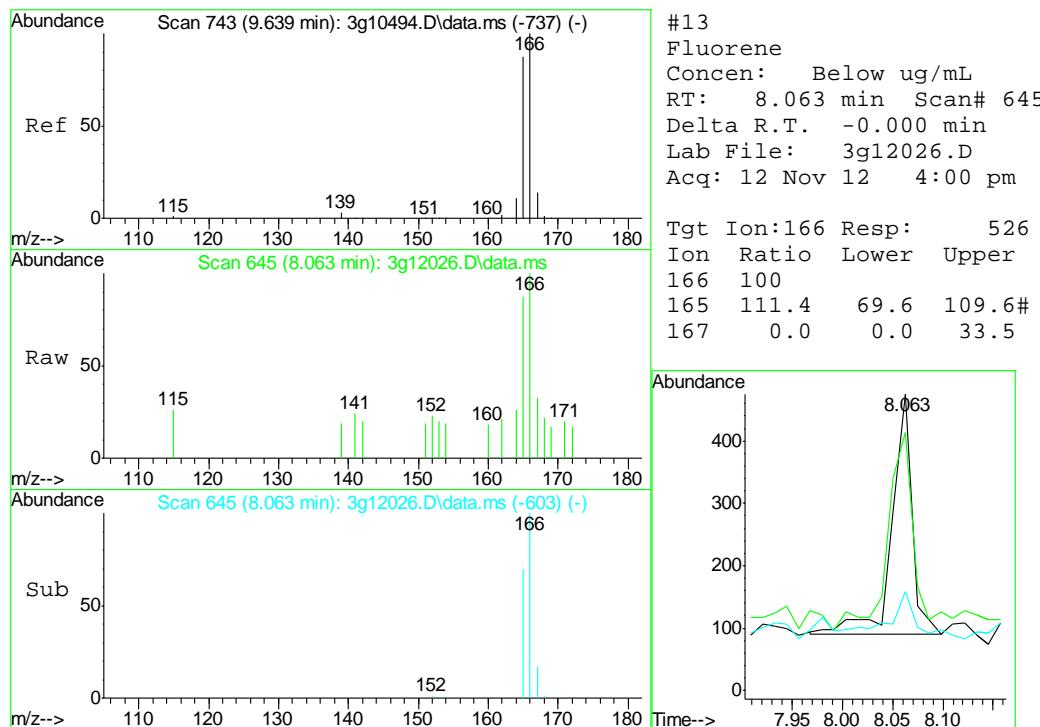


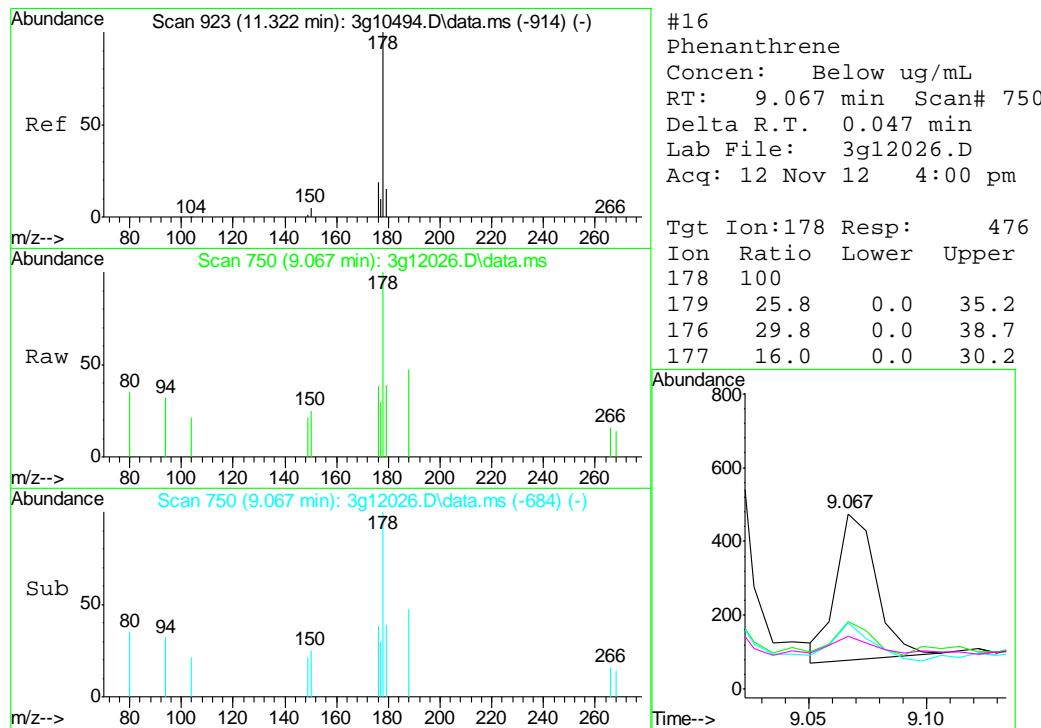
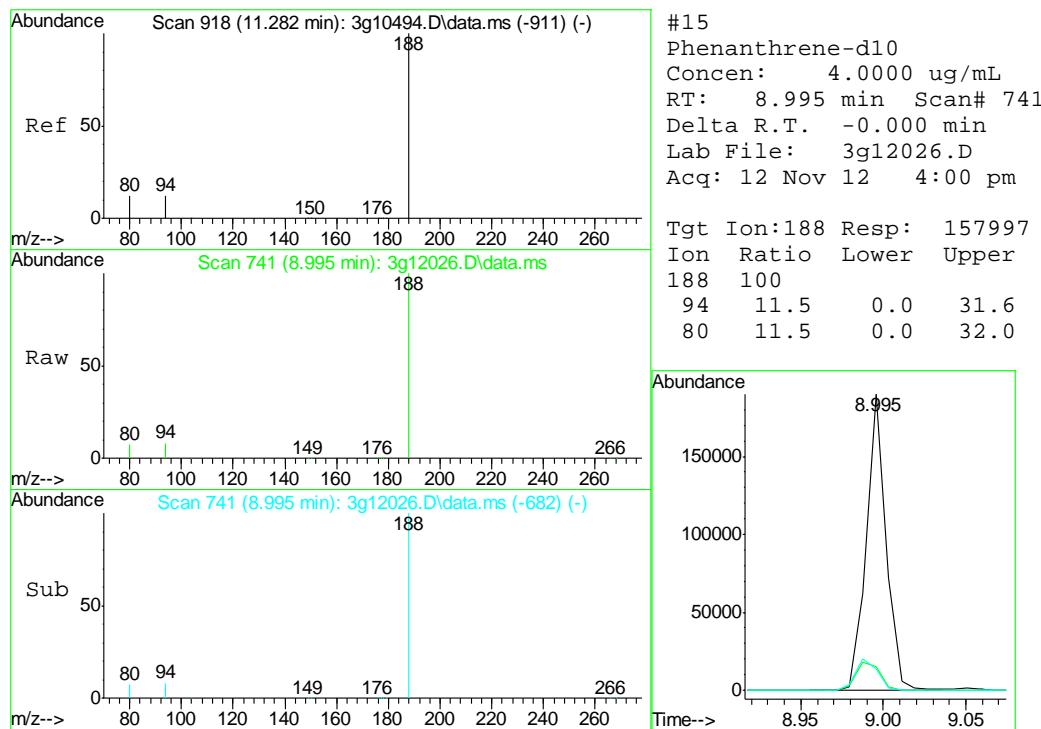


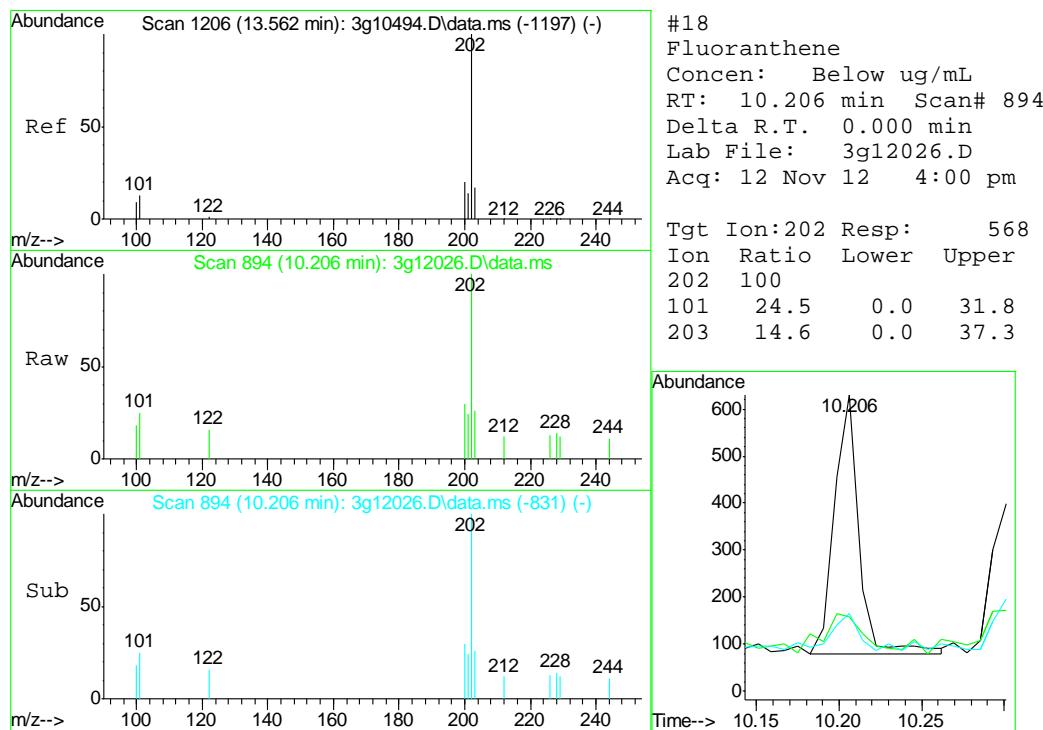
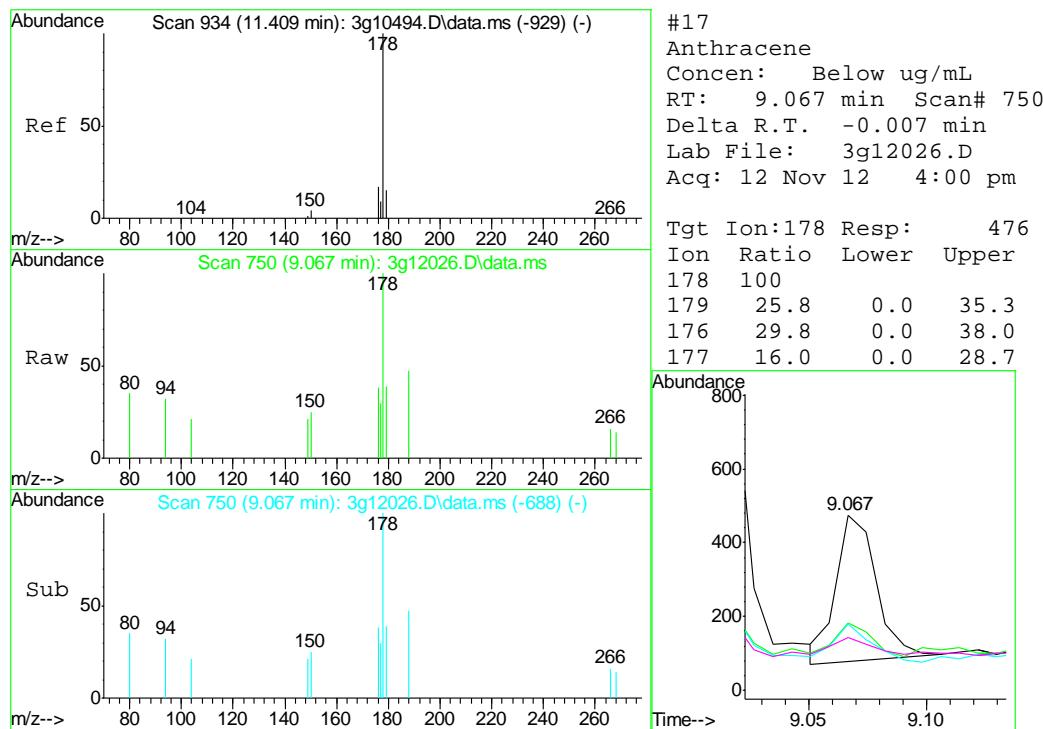


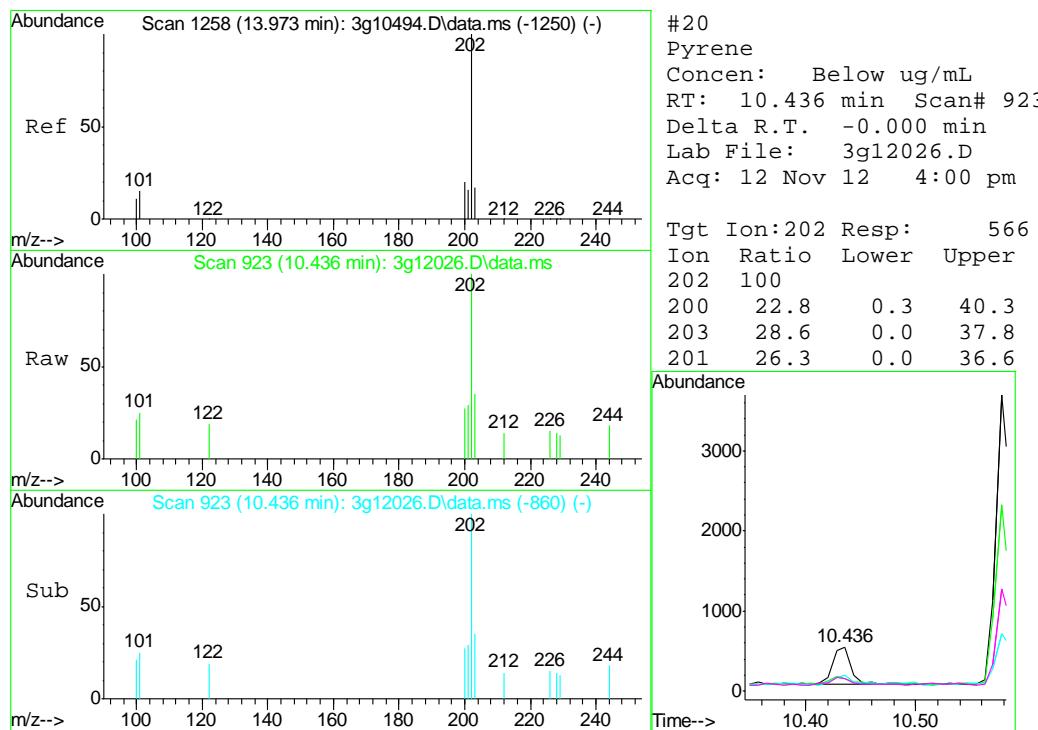
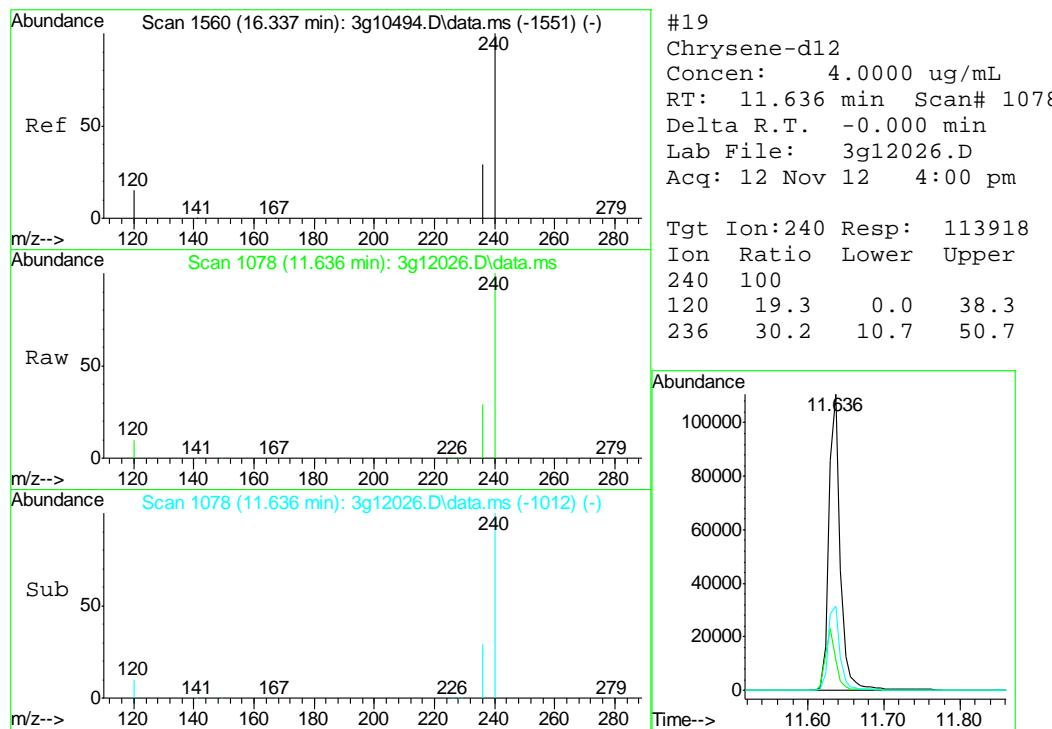


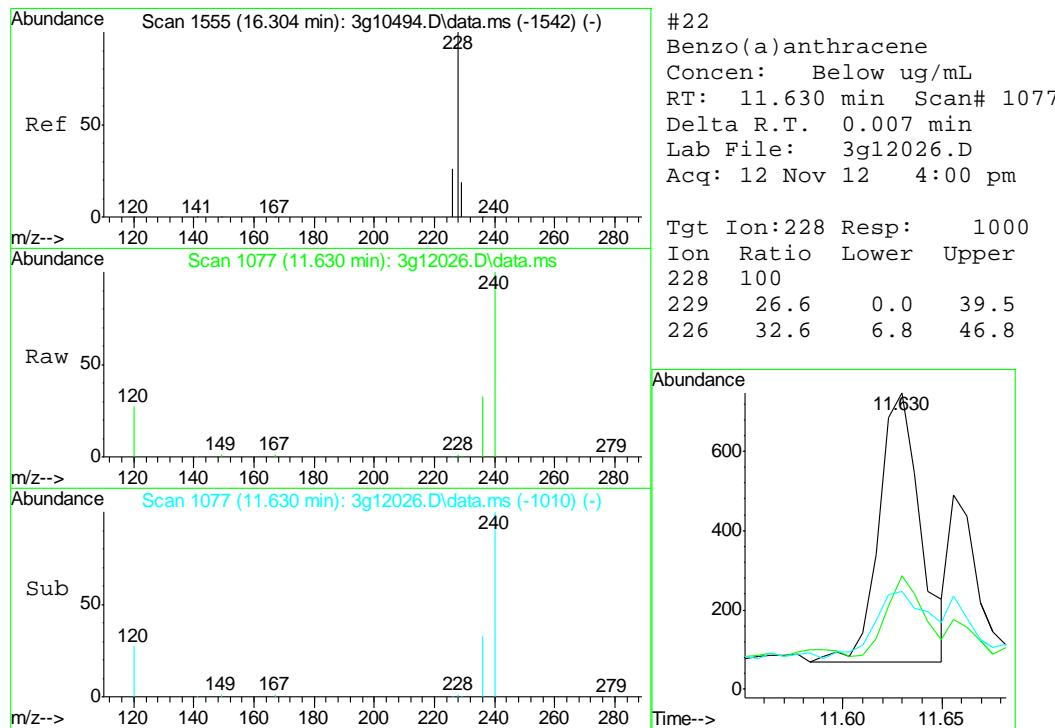
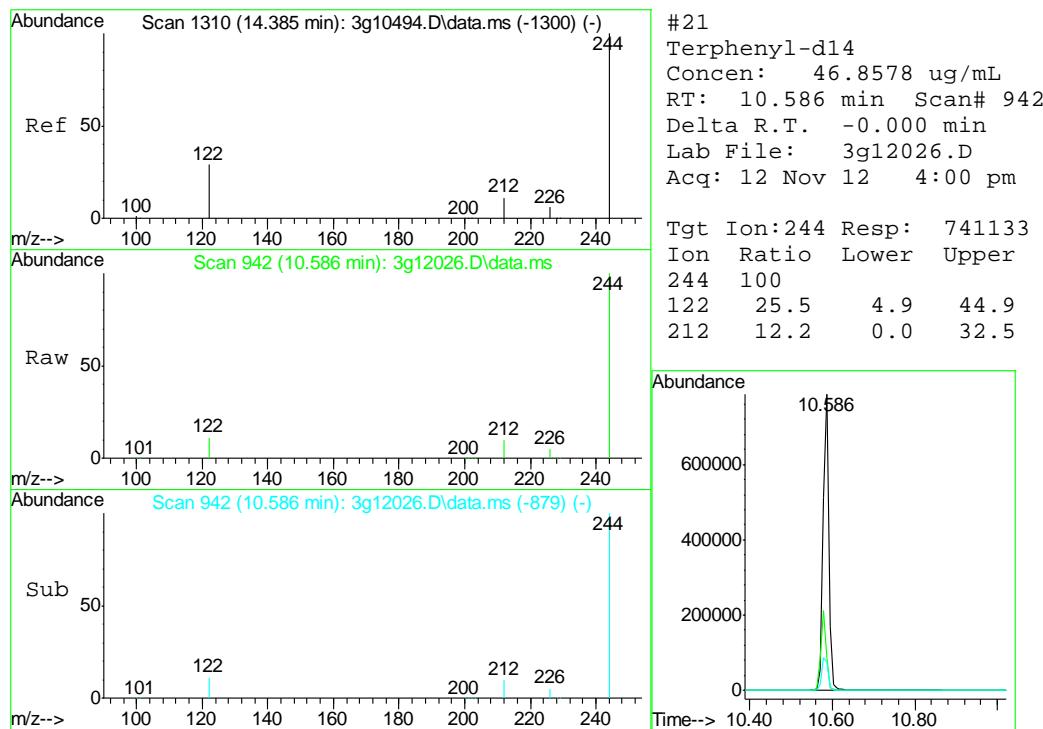


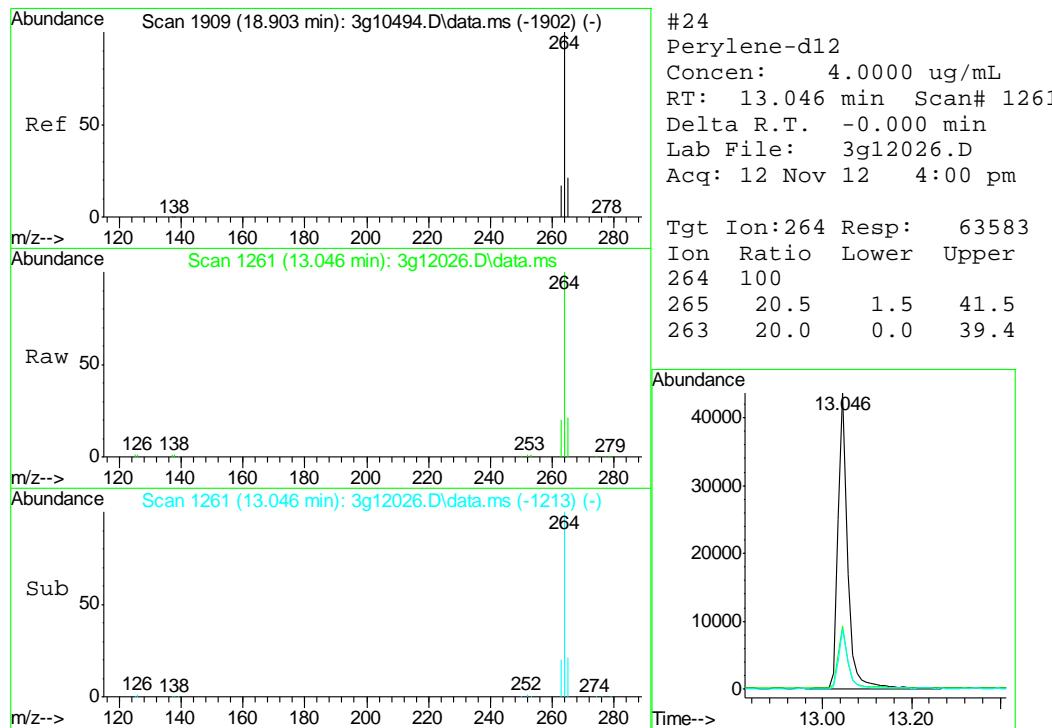
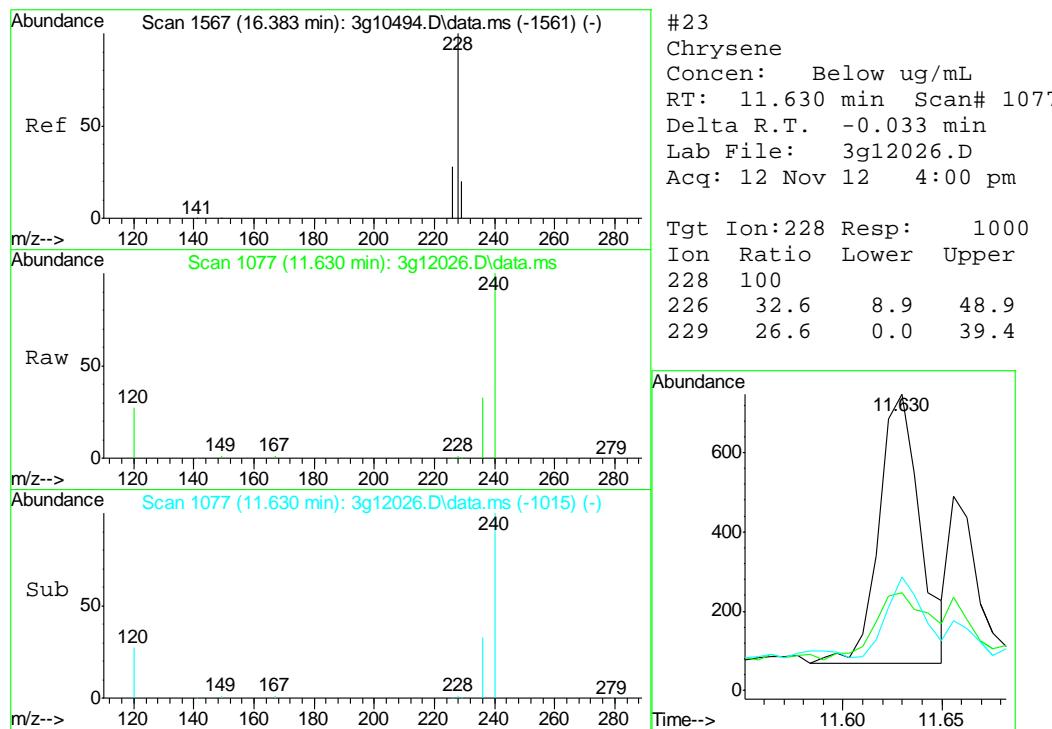


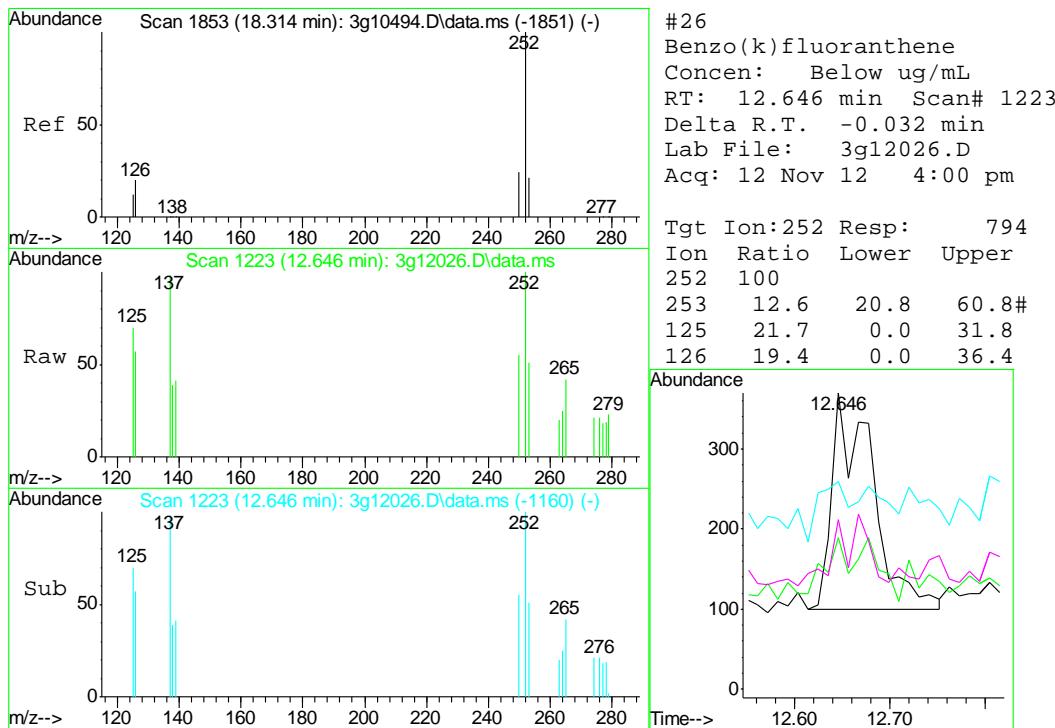
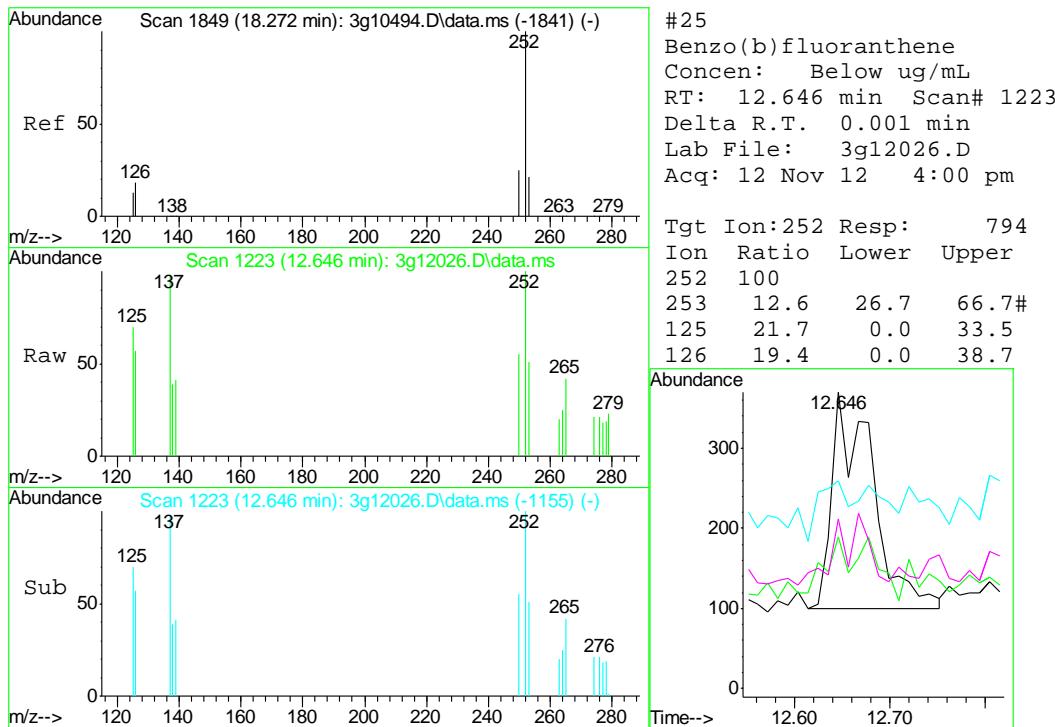


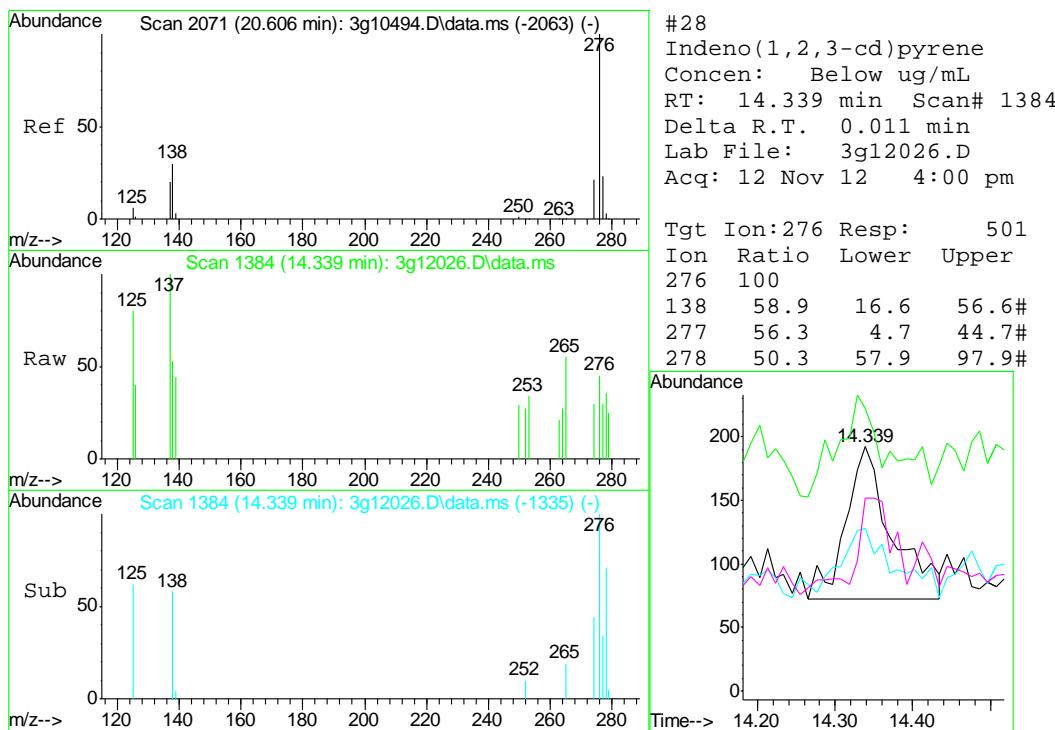
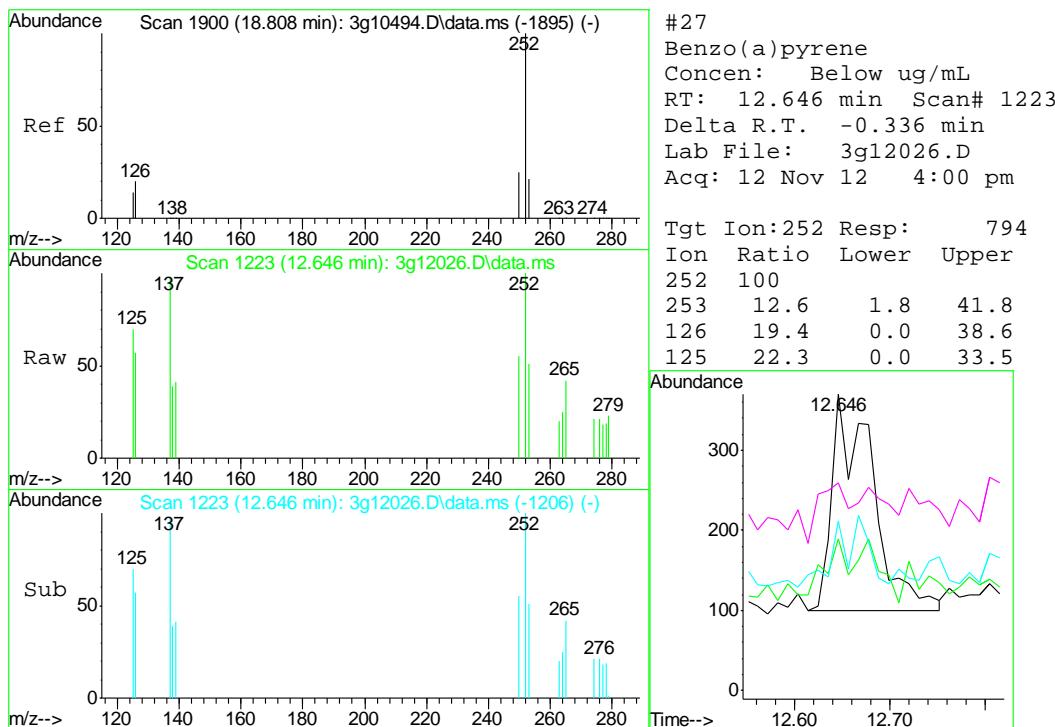


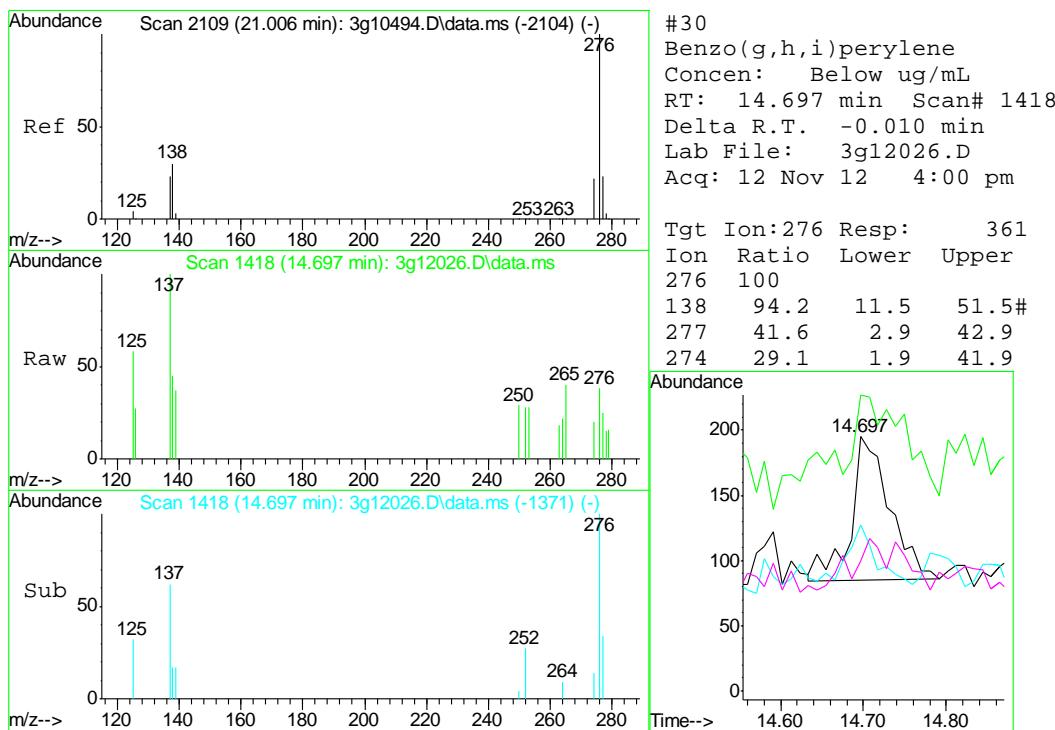
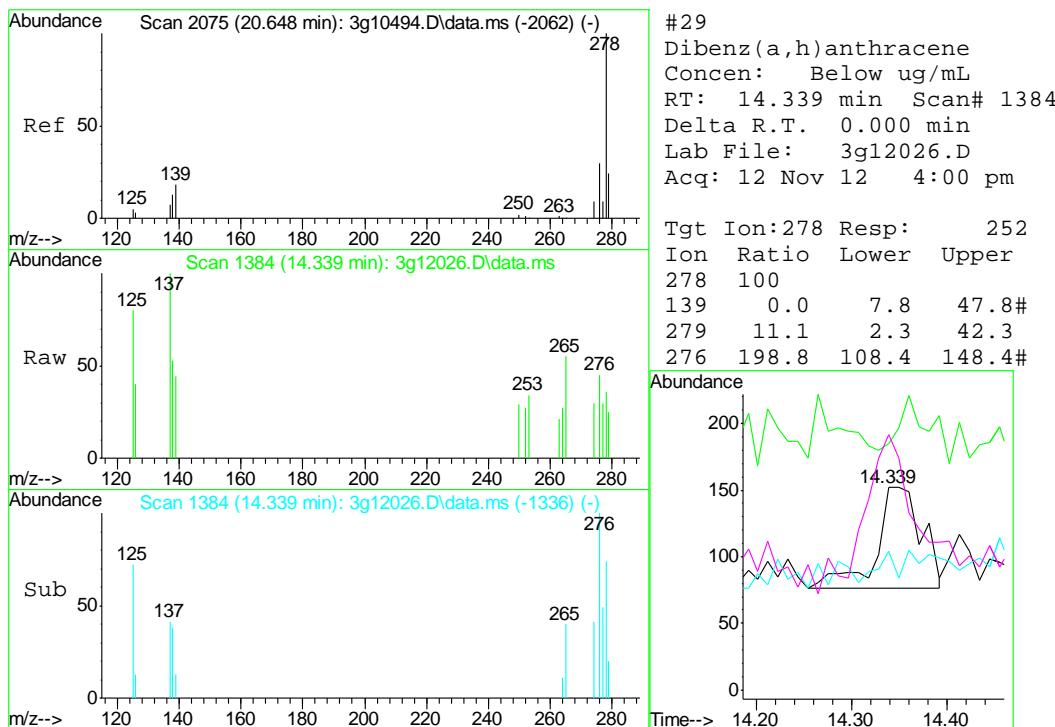














## GC Volatiles

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### QC Data Summaries

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Includes the following where applicable:

- Method Blank Summaries
- Blank Spike Summaries
- Matrix Spike and Duplicate Summaries

**Method Blank Summary**

Job Number: D40798  
 Account: XTOKRWR XTO Energy  
 Project: PCU 296-6A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
GGB1006-MB	GB18409.D	1	11/12/12	SK	n/a	n/a	GGB1006

The QC reported here applies to the following samples:

Method: SW846 8015B

D40798-1

CAS No.	Compound	Result	RL	MDL	Units	Q
	TPH-GRO (C6-C10)	ND	10	5.0	mg/kg	

CAS No.	Surrogate Recoveries	Limits
120-82-1	1,2,4-Trichlorobenzene	91%      60-140%

10.1.1

10

## Blank Spike Summary

Page 1 of 1

Job Number: D40798

Account: XTOKWR XTO Energy

Project: PCU 296-6A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
GGB1006-BS	GB18410.D	1	11/12/12	SK	n/a	n/a	GGB1006

The QC reported here applies to the following samples:

Method: SW846 8015B

D40798-1

CAS No.	Compound	Spike mg/kg	BSP mg/kg	BSP %	Limits
	TPH-GRO (C6-C10)	110	120	109	70-130

CAS No.	Surrogate Recoveries	BSP	Limits
120-82-1	1,2,4-Trichlorobenzene	107%	60-140%

10.2.1

10

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\* = Outside of Control Limits.

# Matrix Spike/Matrix Spike Duplicate Summary

Page 1 of 1

Job Number: D40798

Account: XTOKWR XTO Energy

Project: PCU 296-6A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
D40795-1MS	GB18412.D	1	11/12/12	SK	n/a	n/a	GGB1006
D40795-1MSD	GB18413.D	1	11/12/12	SK	n/a	n/a	GGB1006
D40795-1	GB18411.D	1	11/12/12	SK	n/a	n/a	GGB1006

The QC reported here applies to the following samples:

Method: SW846 8015B

D40798-1

CAS No.	Compound	D40795-1		Spike	MS	MS	MSD	MSD	RPD	Limits Rec/RPD
		mg/kg	Q	mg/kg	mg/kg	%	mg/kg	%		
	TPH-GRO (C6-C10)	ND		149	166	112	164	110	1	70-130/30
<b>10.3.1</b>										
CAS No.	Surrogate Recoveries	MS	MSD	D40795-1			Limits			
120-82-1	1,2,4-Trichlorobenzene	108%	109%	96%	60-140%					

\* = Outside of Control Limits.



## GC Volatiles

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### Raw Data

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## Quantitation Report (QT Reviewed)

Signal #1 : Y:\1\DATA\111212\GB18417.D\FID1A.CH Vial: 11  
 Signal #2 : Y:\1\DATA\111212\GB18417.D\FID2B.CH  
 Acq On : 12 Nov 2012 10:27 pm Operator: StephK  
 Sample : D40798-1, 50X Inst : GC/MS Ins  
 Misc : GC3233,GGB1006,5.019,,100,5,1 Multiplr: 1.00  
 IntFile Signal #1: TVH1.E IntFile Signal #2: FB2.E  
 Quant Time: Nov 13 08:28:18 2012 Quant Results File: TB868GB868SOIL.RES

Quant Method : C:\MSDCHEM\1...\TB868GB868SOIL.M (Chemstation Integrator)  
 Title : 8015B/8021B TVH/BTEX  
 Last Update : Tue Nov 13 08:27:24 2012  
 Response via : Initial Calibration  
 DataAcq Meth : TVB4.M

Volume Inj. :  
 Signal #1 Phase : DB-624 Signal #2 Phase: DB-624  
 Signal #1 Info : 0.53 mm Signal #2 Info : 0.53 mm

Compound	R.T.	Response	Conc	Units
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System Monitoring Compounds

2) S 1,2,4-Trichlorobenzene	14.40	3096218	98.813	%
10) S 1,2,4-Trichlorobenzene (P)	14.40	16404146	100.932	%

Target Compounds

1) H TVH-Gasoline	7.23	2903321	<MDL	mg/L
4) T Methyl-t-butyl-ether	0.00	0	N.D.	ug/L d
5) T Benzene	0.00	0	N.D.	ug/L d
6) T Toluene	7.71	105271	0.266	ug/L
7) T Ethylbenzene	0.00	0	N.D.	ug/L d
8) T m,p-Xylene	0.00	0	N.D.	ug/L d
9) T o-Xylene	0.00	0	N.D.	ug/L d
11) T Naphthalene	14.58	177316	0.899	ug/L

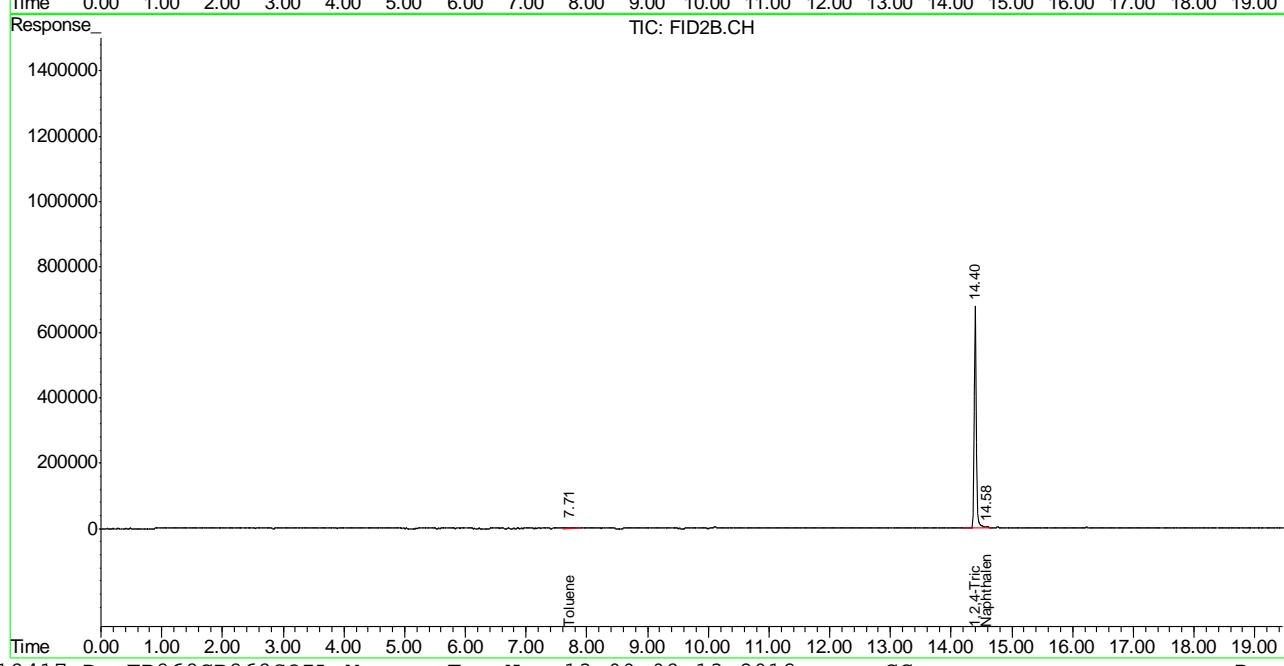
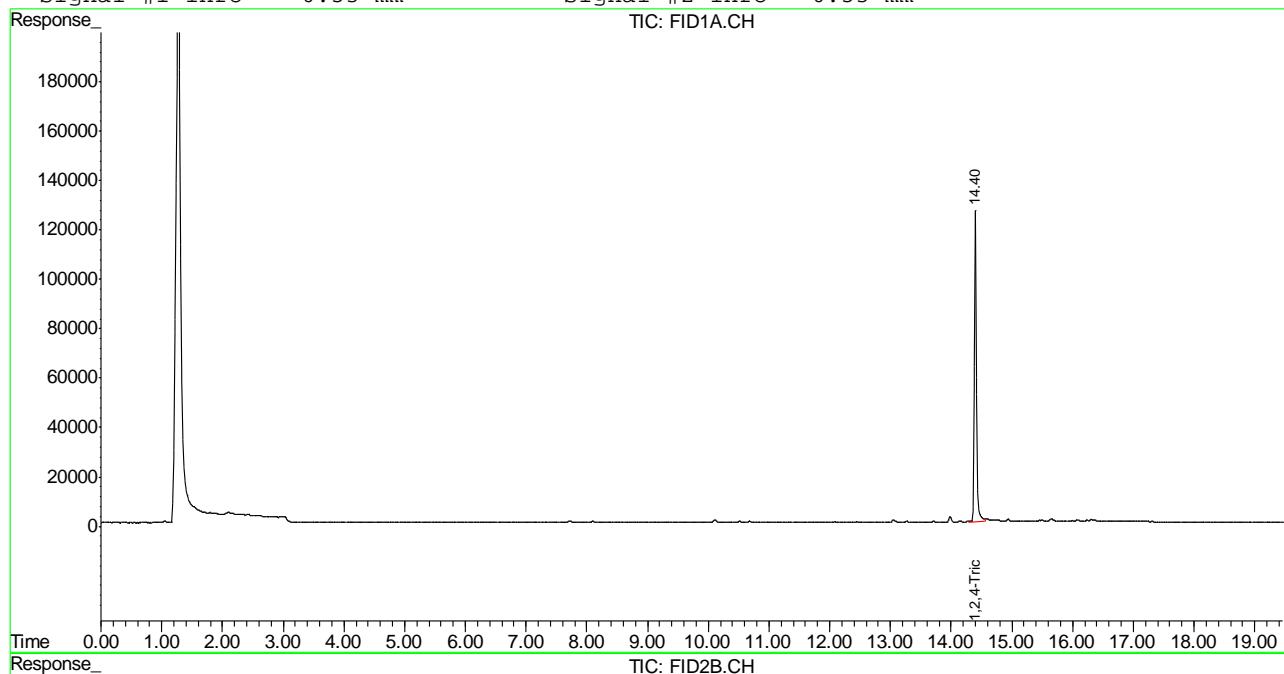
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 (f)=RT Delta > 1/2 Window (m)=manual int.  
 GB18417.D TB868GB868SOIL.M Tue Nov 13 09:08:13 2012 GC

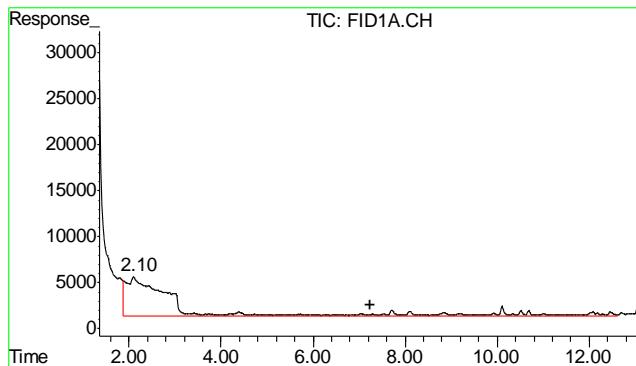
## Quantitation Report (QT Reviewed)

Signal #1 : Y:\1\DATA\111212\GB18417.D\FID1A.CH Vial: 11  
 Signal #2 : Y:\1\DATA\111212\GB18417.D\FID2B.CH  
 Acq On : 12 Nov 2012 10:27 pm Operator: StephK  
 Sample : D40798-1, 50X Inst : GC/MS Ins  
 Misc : GC3233,GGB1006,5.019,,100,5,1 Multiplr: 1.00  
 IntFile Signal #1: TVH1.E IntFile Signal #2: FB2.E  
 Quant Time: Nov 13 9:03 2012 Quant Results File: TB868GB868SOIL.RES

Quant Method : C:\MSDCHEM\1...\TB868GB868SOIL.M (Chemstation Integrator)  
 Title : 8015B/8021B TVH/BTEX  
 Last Update : Tue Nov 13 08:27:24 2012  
 Response via : Multiple Level Calibration  
 DataAcq Meth : TVB4.M

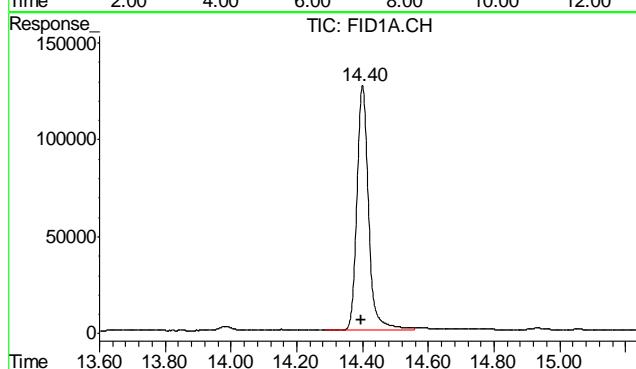
Volume Inj. :  
 Signal #1 Phase : DB-624 Signal #2 Phase: DB-624  
 Signal #1 Info : 0.53 mm Signal #2 Info : 0.53 mm





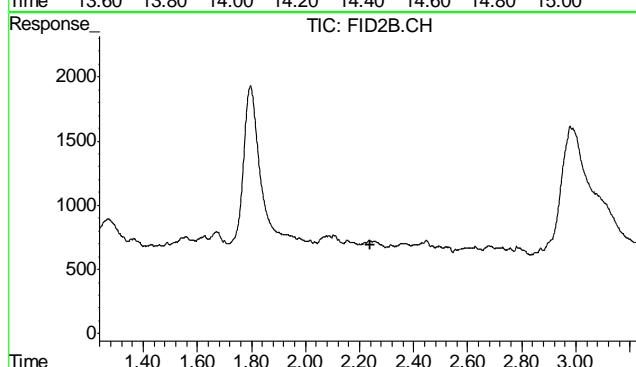
#1 TVH-Gasoline

R.T.: 7.230 min  
Delta R.T.: 0.000 min  
Response: 2903321  
Conc: N.D.



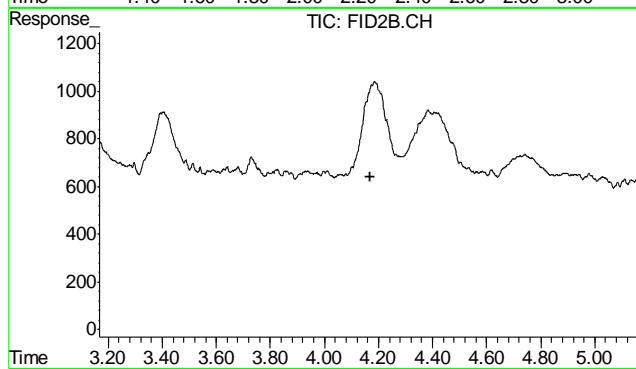
#2 1,2,4-Trichlorobenzene

R.T.: 14.400 min  
Delta R.T.: 0.005 min  
Response: 3096218  
Conc: 98.81 %



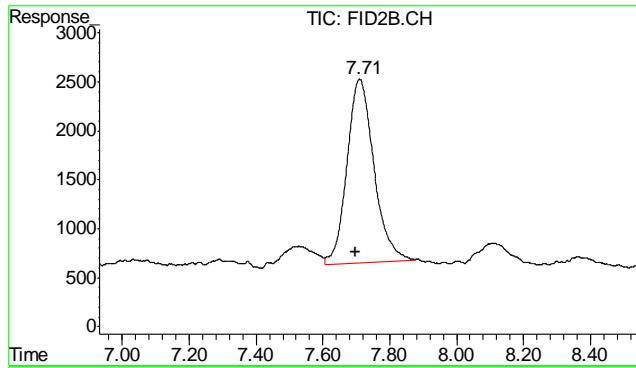
#4 Methyl-t-butyl-ether

R.T.: 0.000 min  
Exp R.T. : 2.238 min  
Response: 0  
Conc: N.D.



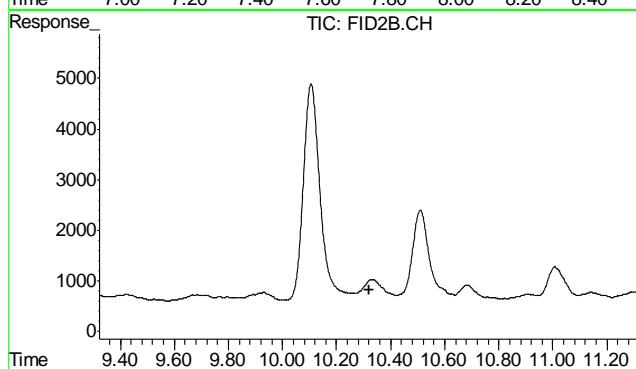
#5 Benzene

R.T.: 0.000 min  
Exp R.T. : 4.168 min  
Response: 0  
Conc: N.D.



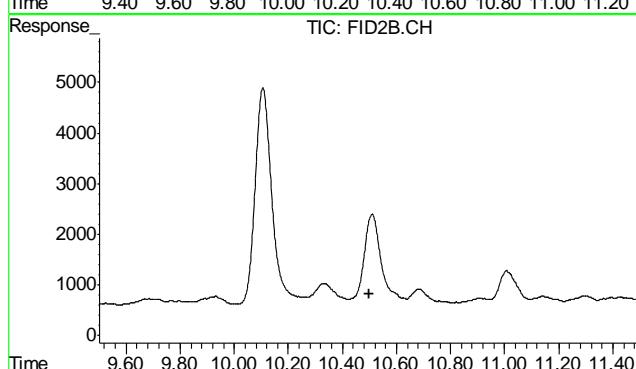
## #6 Toluene

R.T.: 7.710 min  
 Delta R.T.: 0.011 min  
 Response: 105271  
 Conc: 0.27 ug/L



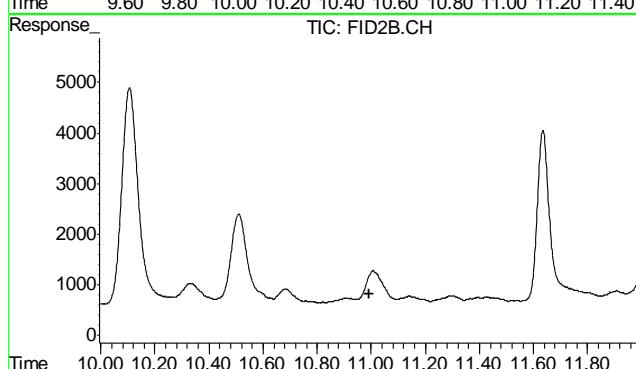
## #7 Ethylbenzene

R.T.: 0.000 min  
 Exp R.T. : 10.322 min  
 Response: 0  
 Conc: N.D.



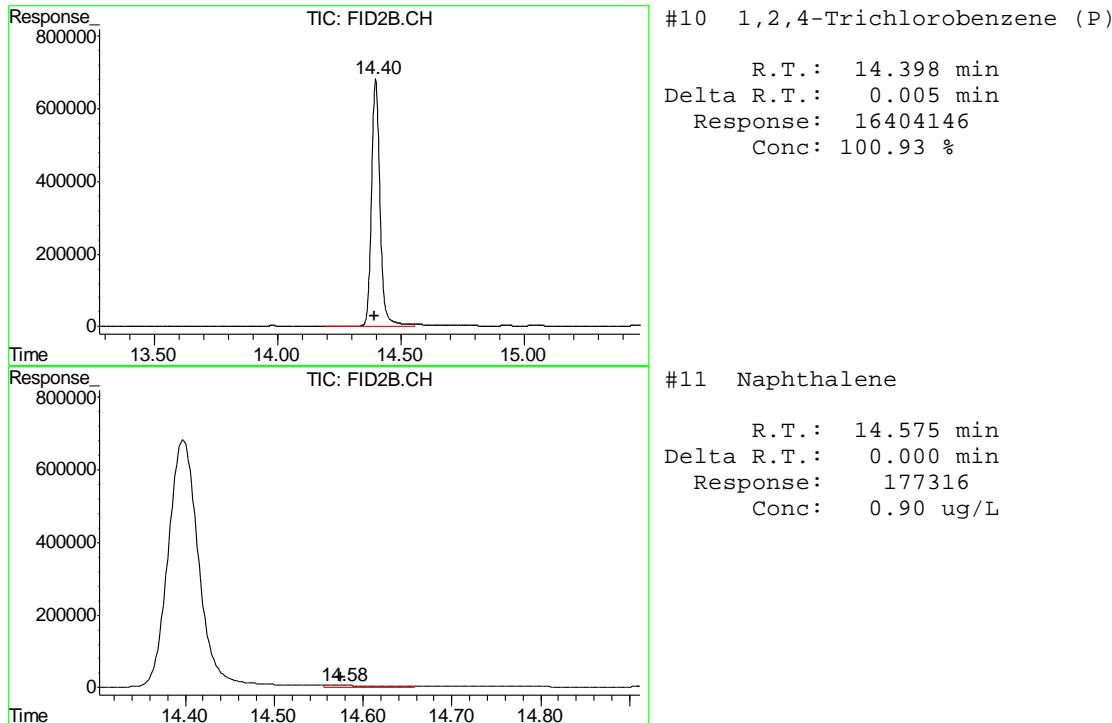
## #8 m,p-Xylene

R.T.: 0.000 min  
 Exp R.T. : 10.500 min  
 Response: 0  
 Conc: N.D.



## #9 o-Xylene

R.T.: 0.000 min  
 Exp R.T. : 10.993 min  
 Response: 0  
 Conc: N.D.



11.1

## Quantitation Report (QT Reviewed)

Signal #1 : Y:\1\DATA\111212\GB18409.D\FID1A.CH Vial: 3  
 Signal #2 : Y:\1\DATA\111212\GB18409.D\FID2B.CH  
 Acq On : 12 Nov 2012 5:43 pm Operator: StephK  
 Sample : MB Inst : GC/MS Ins  
 Misc : GC3233,GGB1006,5.000,,100,5,1 Multiplr: 1.00  
 IntFile Signal #1: TVH1.E IntFile Signal #2: FB2.E  
 Quant Time: Nov 13 08:27:46 2012 Quant Results File: TB868GB868SOIL.RES

Quant Method : C:\MSDCHEM\1...\TB868GB868SOIL.M (Chemstation Integrator)  
 Title : 8015B/8021B TVH/BTEX  
 Last Update : Tue Nov 13 08:27:24 2012  
 Response via : Initial Calibration  
 DataAcq Meth : TVB4.M

Volume Inj. :  
 Signal #1 Phase : DB-624 Signal #2 Phase: DB-624  
 Signal #1 Info : 0.53 mm Signal #2 Info : 0.53 mm

Compound	R.T.	Response	Conc	Units
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System Monitoring Compounds

2) S	1,2,4-Trichlorobenzene	14.39	2846723	90.851 %
10) S	1,2,4-Trichlorobenzene (P)	14.39	15479798	95.244 %

Target Compounds

1) H	TVH-Gasoline	7.23	4545832	<MDL mg/L
4) T	Methyl-t-butyl-ether	0.00	0	N.D. ug/L d
5) T	Benzene	0.00	0	N.D. ug/L d
6) T	Toluene	7.70	107128	0.270 ug/L
7) T	Ethylbenzene	0.00	0	N.D. ug/L d
8) T	m,p-Xylene	0.00	0	N.D. ug/L d
9) T	o-Xylene	0.00	0	N.D. ug/L d
11) T	Naphthalene	14.56	159328	0.808 ug/L

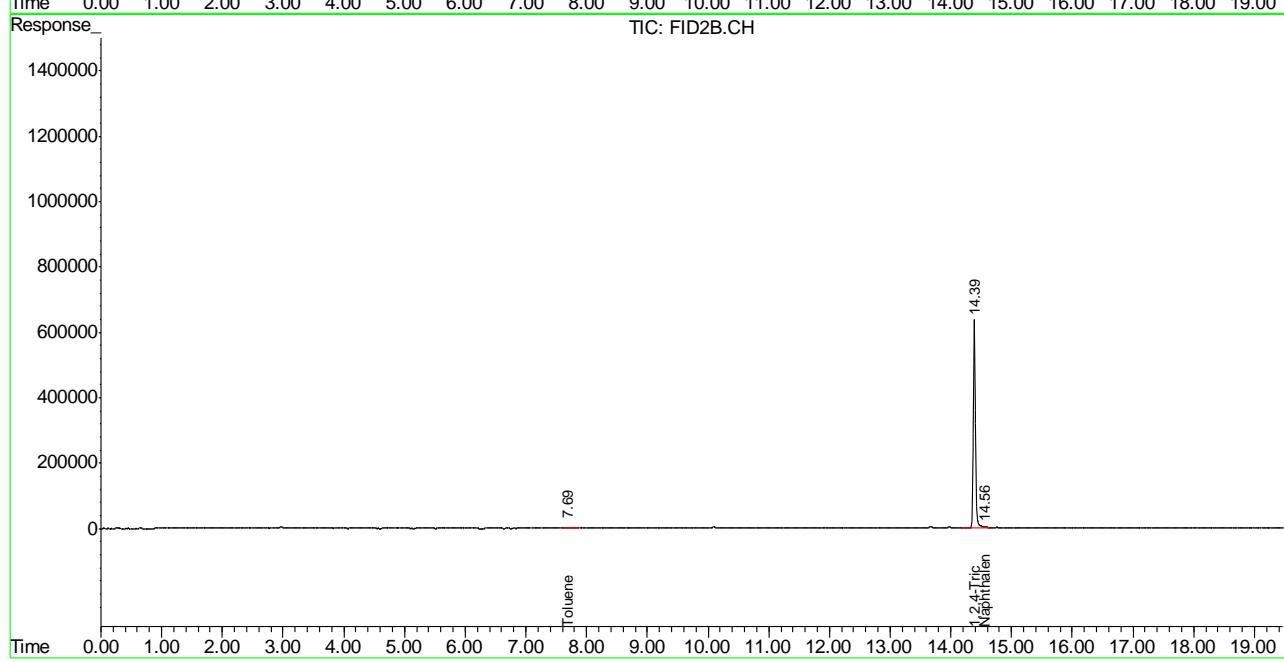
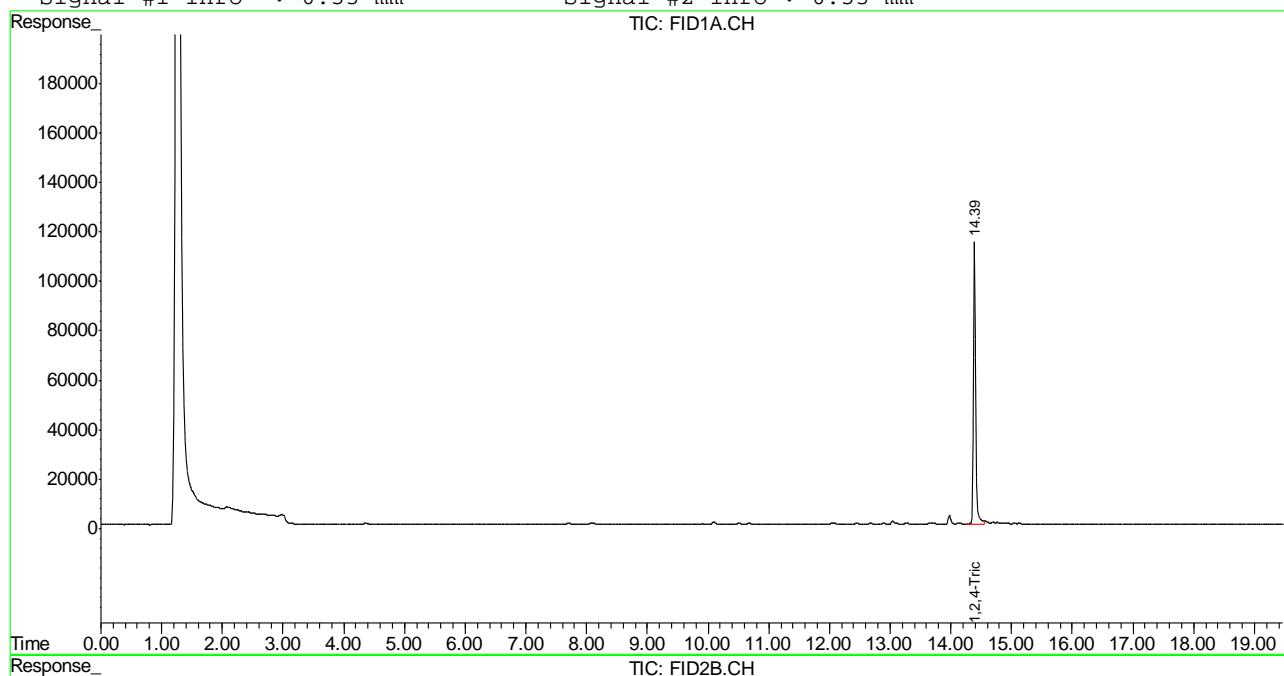
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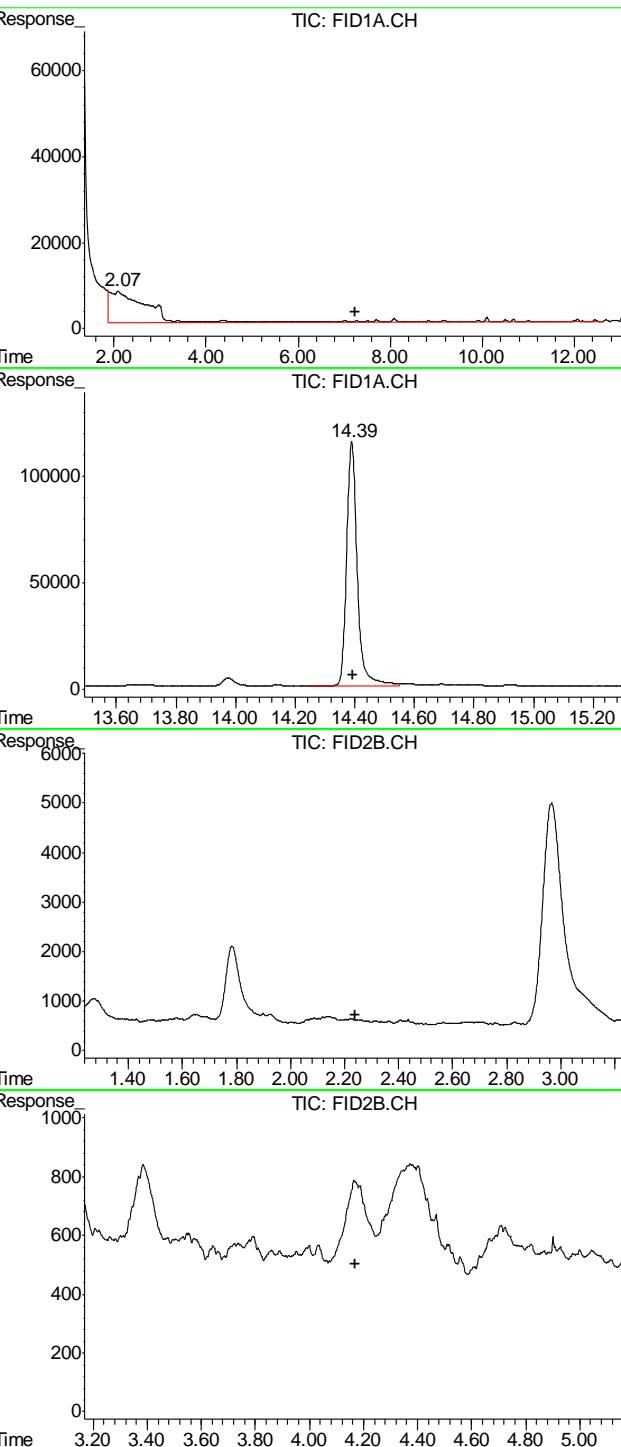
## Quantitation Report (QT Reviewed)

Signal #1 : Y:\1\DATA\111212\GB18409.D\FID1A.CH Vial: 3  
 Signal #2 : Y:\1\DATA\111212\GB18409.D\FID2B.CH  
 Acq On : 12 Nov 2012 5:43 pm Operator: StephK  
 Sample : MB Inst : GC/MS Ins  
 Misc : GC3233,GGB1006,5.000,,100,5,1 Multiplr: 1.00  
 IntFile Signal #1: TVH1.E IntFile Signal #2: FB2.E  
 Quant Time: Nov 13 9:01 2012 Quant Results File: TB868GB868SOIL.RES

Quant Method : C:\MSDCHEM\1...\TB868GB868SOIL.M (Chemstation Integrator)  
 Title : 8015B/8021B TVH/BTEX  
 Last Update : Tue Nov 13 08:27:24 2012  
 Response via : Multiple Level Calibration  
 DataAcq Meth : TVB4.M

Volume Inj. :  
 Signal #1 Phase : DB-624 Signal #2 Phase: DB-624  
 Signal #1 Info : 0.53 mm Signal #2 Info : 0.53 mm



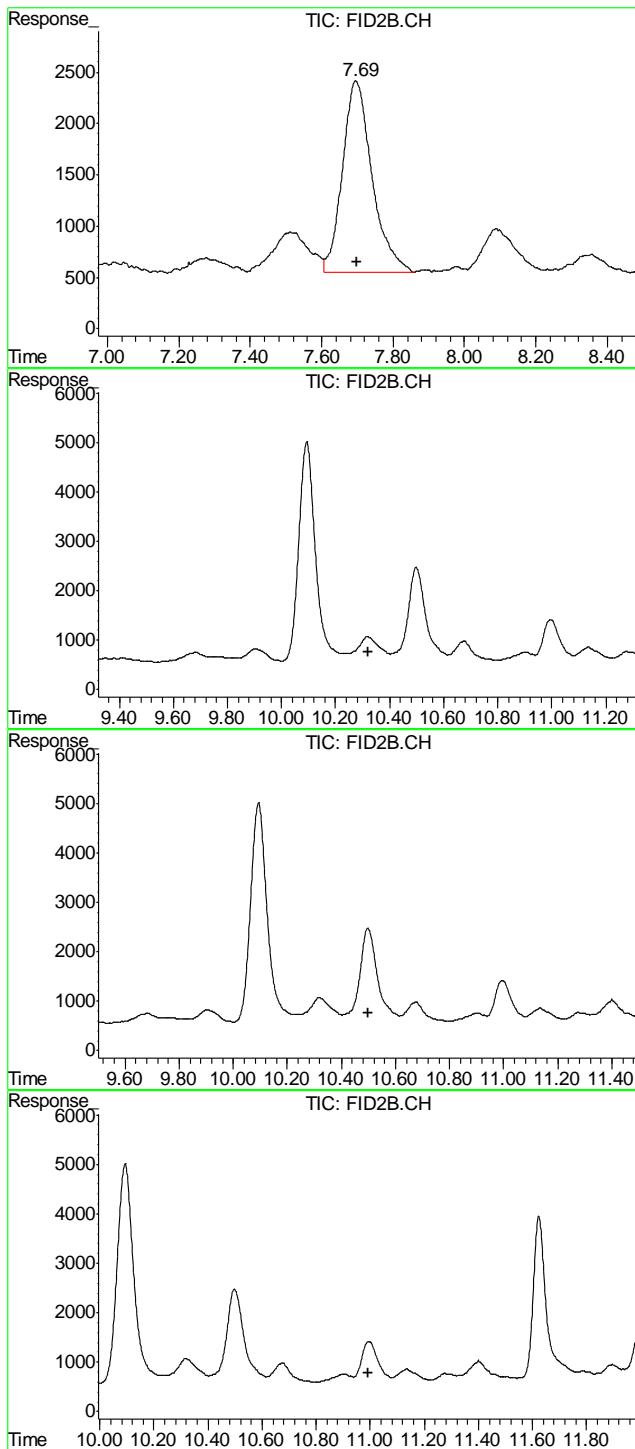


#1 TVH-Gasoline  
 R.T.: 7.230 min  
 Delta R.T.: 0.000 min  
 Response: 4545832  
 Conc: N.D.

#2 1,2,4-Trichlorobenzene  
 R.T.: 14.390 min  
 Delta R.T.: -0.005 min  
 Response: 2846723  
 Conc: 90.85 %

#4 Methyl-t-butyl-ether  
 R.T.: 0.000 min  
 Exp R.T. : 2.238 min  
 Response: 0  
 Conc: N.D.

#5 Benzene  
 R.T.: 0.000 min  
 Exp R.T. : 4.168 min  
 Response: 0  
 Conc: N.D.



#6 Toluene

R.T.: 7.696 min  
Delta R.T.: -0.003 min  
Response: 107128  
Conc: 0.27 ug/L

#7 Ethylbenzene

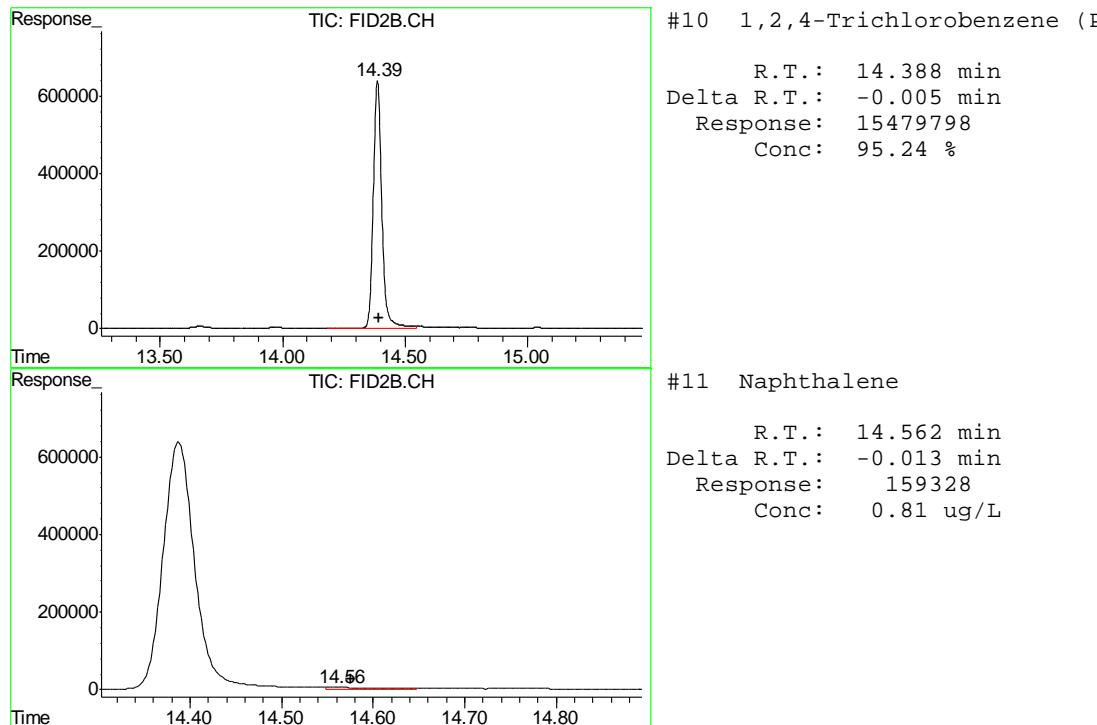
R.T.: 0.000 min  
Exp R.T. : 10.322 min  
Response: 0  
Conc: N.D.

#8 m,p-Xylene

R.T.: 0.000 min  
Exp R.T. : 10.500 min  
Response: 0  
Conc: N.D.

#9 o-Xylene

R.T.: 0.000 min  
Exp R.T. : 10.993 min  
Response: 0  
Conc: N.D.



11.2.1

11



## GC Semi-volatiles

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### QC Data Summaries

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Includes the following where applicable:

- Method Blank Summaries
- Blank Spike Summaries
- Matrix Spike and Duplicate Summaries

**Method Blank Summary**

Job Number: D40798  
 Account: XTOKRWR XTO Energy  
 Project: PCU 296-6A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
OP6948-MB	FD19400.D	1	11/12/12	AV	11/12/12	OP6948	GFD976

The QC reported here applies to the following samples:

Method: SW846-8015B

D40798-1

CAS No.	Compound	Result	RL	MDL	Units	Q
	TPH-DRO (C10-C28)	ND	13	8.7	mg/kg	

CAS No.	Surrogate Recoveries	Limits
84-15-1	o-Terphenyl	101% 35-130%

## Blank Spike Summary

Page 1 of 1

Job Number: D40798

Account: XTOKRWR XTO Energy

Project: PCU 296-6A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
OP6948-BS	FD19402.D	1	11/12/12	AV	11/12/12	OP6948	GFD976

The QC reported here applies to the following samples:

Method: SW846-8015B

D40798-1

CAS No.	Compound	Spike mg/kg	BSP mg/kg	BSP %	Limits
	TPH-DRO (C10-C28)	667	630	94	48-130

CAS No.	Surrogate Recoveries	BSP	Limits
84-15-1	o-Terphenyl	109%	35-130%

\* = Outside of Control Limits.

12.2.1

12

# Matrix Spike/Matrix Spike Duplicate Summary

Page 1 of 1

Job Number: D40798

Account: XTOKWR XTO Energy

Project: PCU 296-6A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
OP6948-MS	FD19404.D	1	11/12/12	AV	11/12/12	OP6948	GFD976
OP6948-MSD	FD19406.D	1	11/12/12	AV	11/12/12	OP6948	GFD976
D40799-1	FD19430.D	1	11/12/12	AV	11/12/12	OP6948	GFD976

The QC reported here applies to the following samples:

Method: SW846-8015B

D40798-1

CAS No.	Compound	D40799-1		Spike	MS	MS	MSD	MSD	RPD	Limits Rec/RPD
		mg/kg	Q	mg/kg	mg/kg	%	mg/kg	%		
	TPH-DRO (C10-C28)	23.1		720	608	81	612	82	1	20-168/30
<hr/>										
CAS No.	Surrogate Recoveries	MS		MSD		D40799-1	Limits			
84-15-1	o-Terphenyl	94%		92%		90%	35-130%			

\* = Outside of Control Limits.

12.3.1  
12



## GC Semi-volatiles

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### Raw Data

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## Quantitation Report (QT Reviewed)

Data File : C:\MSDCHEM\2\DATA\2012\NOV\FD111212.SEC\FD19426.D Vial: 66  
 Acq On : 11-12-2012 08:23:32 PM Operator: ashleyv  
 Sample : D40798-1 Inst : FID5  
 Misc : OP6948,GFD976,30.06,,,2,1 Multiplr: 1.00  
 IntFile : autoint1.e  
 Quant Time: Nov 13 09:26:06 2012 Quant Results File: DRO-GFD939R.RES

Quant Method : C:\MSDCHEM\2...\DRO-GFD939R.M (Chemstation Integrator)  
 Title : 8015B TEH  
 Last Update : Thu Nov 01 15:20:48 2012  
 Response via : Initial Calibration  
 DataAcq Meth : DRODUAL.M

Volume Inj. : 1ul  
 Signal Phase : RTX-5  
 Signal Info : 530um

Compound	R.T.	Response	Conc Units
<hr/>			
System Monitoring Compounds			
1) S O-Terphenyl	9.08	36979138	911.475 mg/L
<hr/>			
Target Compounds			
2) H TPH-DRO (c10-c28)	7.04	2519502	69.193 mg/L

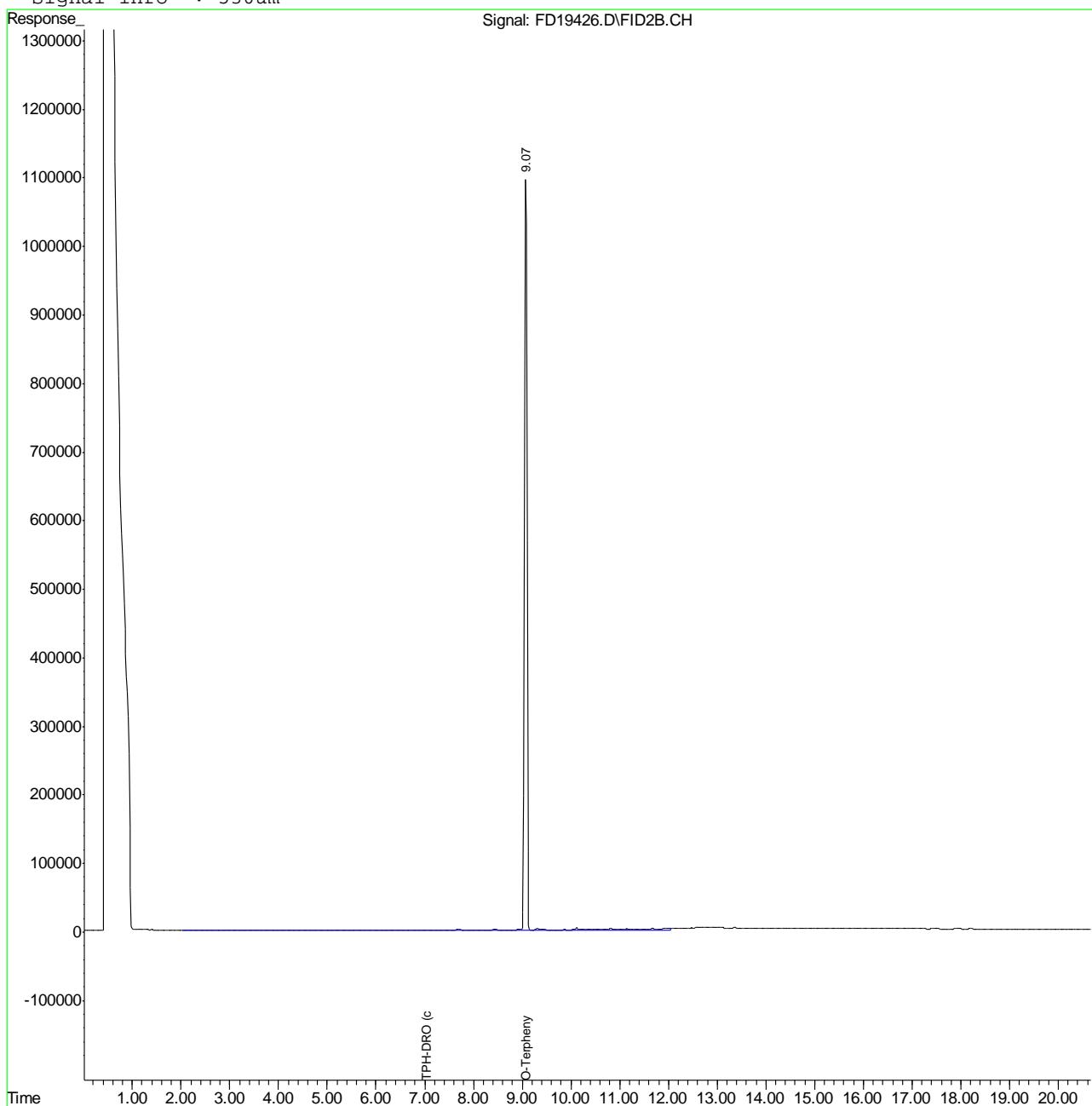
(f)=RT Delta > 1/2 Window (m)=manual int.  
 FD19426.D DRO-GFD939R.M Tue Nov 13 09:40:39 2012 GC

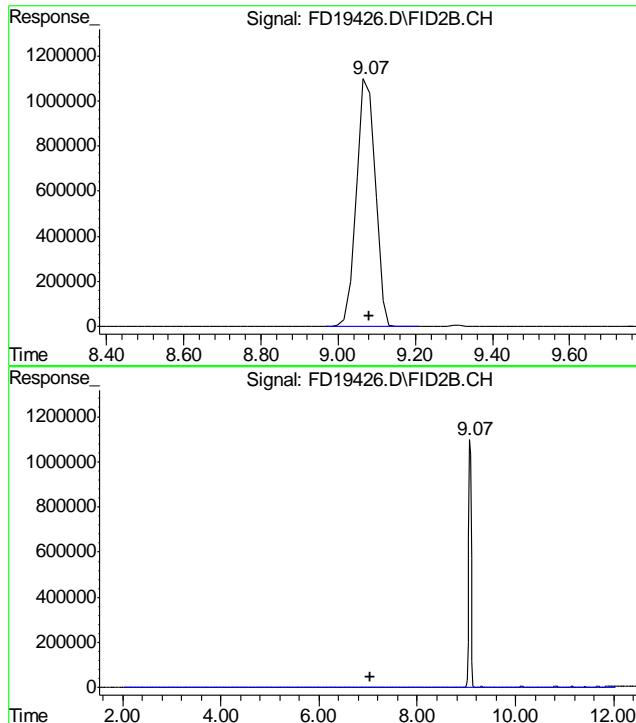
## Quantitation Report (QT Reviewed)

Data File : C:\MSDCHEM\2\DATA\2012\NOV\FD111212.SEC\FD19426.D Vial: 66  
 Acq On : 11-12-2012 08:23:32 PM Operator: ashleyv  
 Sample : D40798-1 Inst : FID5  
 Misc : OP6948,GFD976,30.06,,,2,1 Multiplr: 1.00  
 IntFile : autoint1.e  
 Quant Time: Nov 13 9:26 2012 Quant Results File: DRO-GFD939R.RES

Quant Method : C:\MSDCHEM\2...\DRO-GFD939R.M (Chemstation Integrator)  
 Title : 8015B TEH  
 Last Update : Thu Nov 01 15:20:48 2012  
 Response via : Multiple Level Calibration  
 DataAcq Meth : DRODUAL.M

Volume Inj. : 1ul  
 Signal Phase : RTX-5  
 Signal Info : 530um





#1 O-Terphenyl  
 R.T.: 9.079 min  
 Delta R.T.: -0.001 min  
 Response: 36979138  
 Conc: 911.48 mg/L

#2 TPH-DRO (c10-c28)  
 R.T.: 7.035 min  
 Delta R.T.: 0.000 min  
 Response: 2519502  
 Conc: 69.19 mg/L

## Quantitation Report (QT Reviewed)

Data File : C:\MSDCHEM\2\DATA\2012\NOV\FD111212.SEC\FD19400.D Vial: 53  
 Acq On : 11-12-2012 02:28:54 PM Operator: ashleyv  
 Sample : OP6948-MB Inst : FID5  
 Misc : OP6948,GFD976,30.00,,,2,1 Multiplr: 1.00  
 IntFile : autoint1.e  
 Quant Time: Nov 13 09:25:53 2012 Quant Results File: DRO-GFD939R.RES

Quant Method : C:\MSDCHEM\2...\DRO-GFD939R.M (Chemstation Integrator)  
 Title : 8015B TEH  
 Last Update : Thu Nov 01 15:20:48 2012  
 Response via : Initial Calibration  
 DataAcq Meth : DRODUAL.M

Volume Inj. : 1ul  
 Signal Phase : RTX-5  
 Signal Info : 530um

Compound	R.T.	Response	Conc Units
<hr/>			
System Monitoring Compounds			
1) S O-Terphenyl	9.08	40924247	1008.716 mg/L
<hr/>			
Target Compounds			
2) H TPH-DRO (c10-c28)	7.04	931533	25.583 mg/L

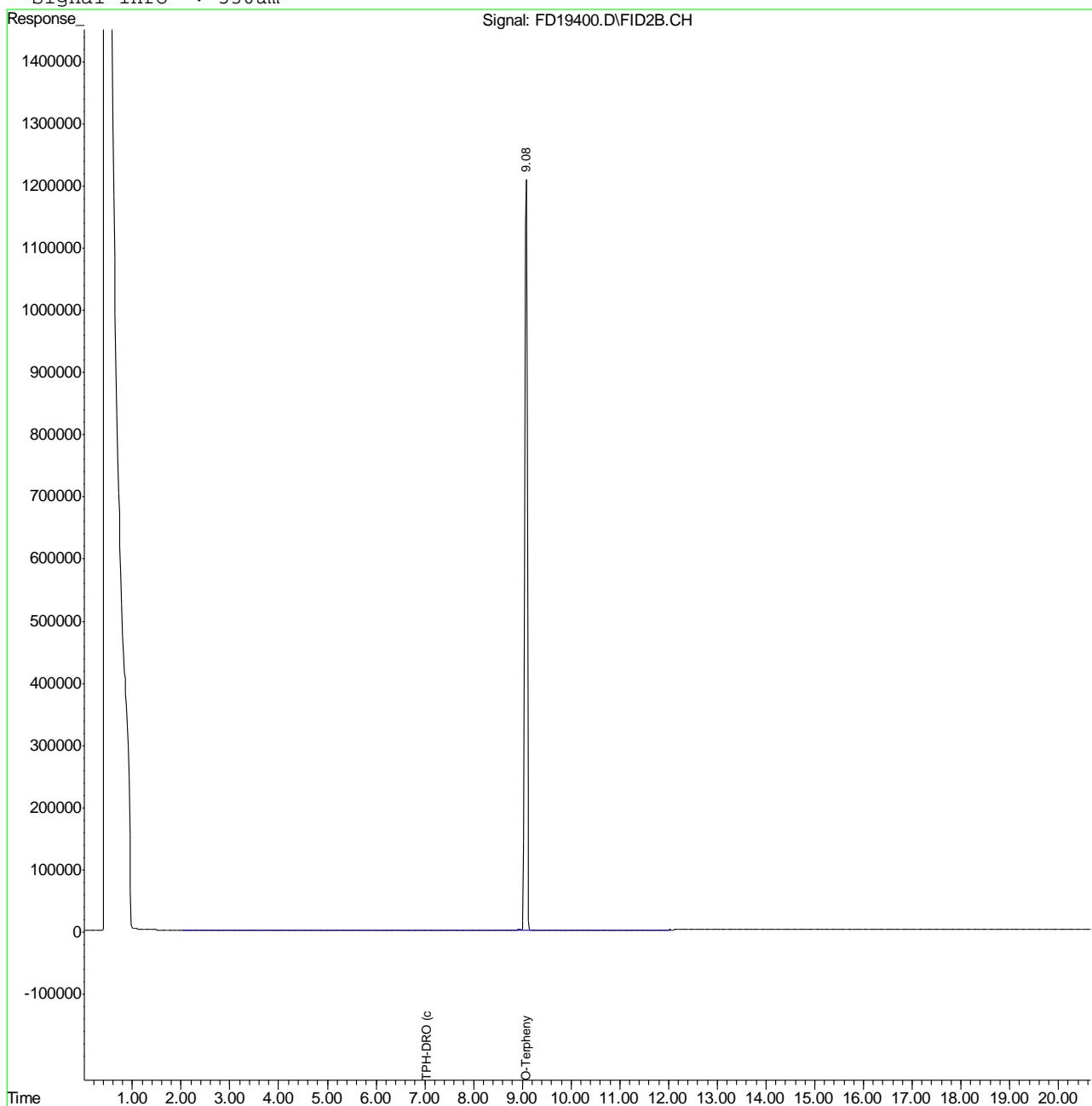
(f)=RT Delta > 1/2 Window (m)=manual int.  
 FD19400.D DRO-GFD939R.M Tue Nov 13 09:40:26 2012 GC

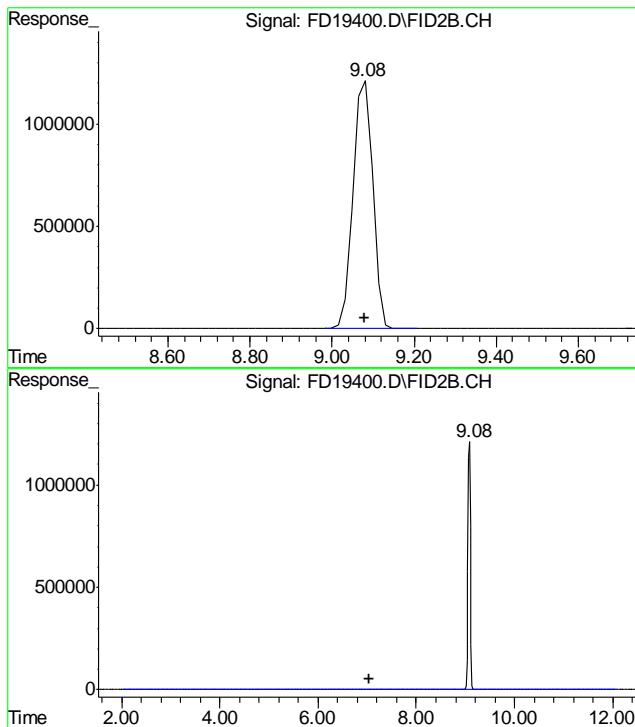
## Quantitation Report (QT Reviewed)

Data File : C:\MSDCHEM\2\DATA\2012\NOV\FD111212.SEC\FD19400.D Vial: 53  
 Acq On : 11-12-2012 02:28:54 PM Operator: ashleyv  
 Sample : OP6948-MB Inst : FID5  
 Misc : OP6948,GFD976,30.00,,,2,1 Multiplr: 1.00  
 IntFile : autoint1.e  
 Quant Time: Nov 13 9:25 2012 Quant Results File: DRO-GFD939R.RES

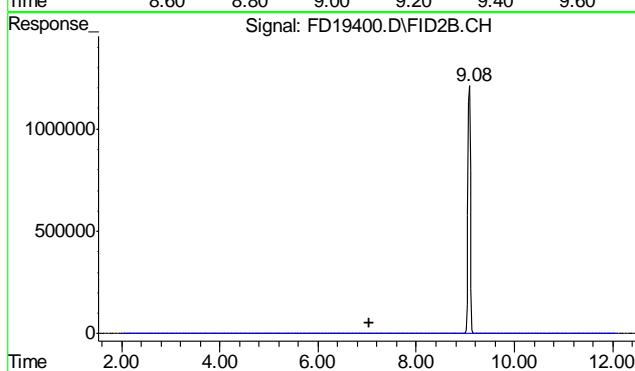
Quant Method : C:\MSDCHEM\2...\DRO-GFD939R.M (Chemstation Integrator)  
 Title : 8015B TEH  
 Last Update : Thu Nov 01 15:20:48 2012  
 Response via : Multiple Level Calibration  
 DataAcq Meth : DRODUAL.M

Volume Inj. : 1uL  
 Signal Phase : RTX-5  
 Signal Info : 530um





#1 O-Terphenyl  
R.T.: 9.082 min  
Delta R.T.: 0.002 min  
Response: 40924247  
Conc: 1008.72 mg/L



#2 TPH-DRO (c10-c28)  
R.T.: 7.035 min  
Delta R.T.: 0.000 min  
Response: 931533  
Conc: 25.58 mg/L m



## Metals Analysis

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### QC Data Summaries

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Includes the following where applicable:

- Method Blank Summaries
- Matrix Spike and Duplicate Summaries
- Blank Spike and Lab Control Sample Summaries
- Serial Dilution Summaries

BLANK RESULTS SUMMARY  
Part 2 - Method Blanks

Login Number: D40798  
Account: XTOKRWR - XTO Energy  
Project: PCU 296-6A

QC Batch ID: MP8869  
Matrix Type: SOLID

Methods: SW846 6010C  
Units: mg/kg

Prep Date:

11/13/12

Metal	RL	IDL	MDL	MB raw	final
Aluminum	10	.96	.57		
Antimony	3.0	.17	.12		
Arsenic	2.5	.44	.56		
Barium	1.0	.01	.11	0.47	<1.0
Beryllium	1.0	.13	.15		
Boron	5.0	.1	.06		
Cadmium	1.0	.06	.036	0.010	<1.0
Calcium	40	.54	9		
Chromium	1.0	.03	.03	0.060	<1.0
Cobalt	0.50	.04	.07		
Copper	1.0	.12	.15	-0.14	<1.0
Iron	7.0	.12	.87		
Lead	5.0	.19	.24	0.080	<5.0
Lithium	0.20	.05	.054		
Magnesium	20	.65	.98		
Manganese	0.50	.12	.022		
Molybdenum	1.0	.21	.08		
Nickel	3.0	.05	.026	0.12	<3.0
Phosphorus	10	1.4	1.9		
Potassium	200	6.1	7		
Selenium	5.0	.48	.36	-0.030	<5.0
Silicon	5.0	.29	.37		
Silver	3.0	.04	.06	-0.050	<3.0
Sodium	40	.59	1.9		
Strontium	5.0	.004	.017		
Thallium	1.0	.29	.53		
Tin	5.0	1.2	2		
Titanium	1.0	.01	.038		
Uranium	5.0	.22	.26		
Vanadium	1.0	.02	.036		
Zinc	3.0	.05	.37	0.27	<3.0

Associated samples MP8869: D40798-1

Results < IDL are shown as zero for calculation purposes  
(\*) Outside of QC limits

BLANK RESULTS SUMMARY  
Part 2 - Method Blanks

Login Number: D40798  
Account: XTOKRWR - XTO Energy  
Project: PCU 296-6A

QC Batch ID: MP8869  
Matrix Type: SOLID

Methods: SW846 6010C  
Units: mg/kg

Prep Date:

Metal

(anr) Analyte not requested

## MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D40798  
 Account: XTOKRWR - XTO Energy  
 Project: PCU 296-6A

QC Batch ID: MP8869  
 Matrix Type: SOLID

Methods: SW846 6010C  
 Units: mg/kg

Prep Date:

11/13/12

Metal	D40778-1 Original MS	Spikelot ICPALL2	% Rec	QC Limits
Aluminum				
Antimony				
Arsenic				
Barium	3150	3380	233	98.8 75-125
Beryllium				
Boron	anr			
Cadmium	0.0	51.1	58.2	87.8 75-125
Calcium				
Chromium	65.8	117	58.2	88.0 75-125
Cobalt				
Copper	14.3	68.1	58.2	92.5 75-125
Iron				
Lead	10.3	110	116	85.7 75-125
Lithium				
Magnesium				
Manganese				
Molybdenum				
Nickel	23.9	71.5	58.2	81.8 75-125
Phosphorus				
Potassium				
Selenium	0.64	104	116	88.8 75-125
Silicon				
Silver	0.13	21.4	23.3	91.4 75-125
Sodium				
Strontium				
Thallium				
Tin				
Titanium				
Uranium				
Vanadium				
Zinc	52.2	104	58.2	89.0 75-125

Associated samples MP8869: D40798-1

Results < IDL are shown as zero for calculation purposes  
 (\*) Outside of QC limits

MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D40798  
Account: XTOKRWR - XTO Energy  
Project: PCU 296-6A

QC Batch ID: MP8869  
Matrix Type: SOLID

Methods: SW846 6010C  
Units: mg/kg

Prep Date:

Metal

(N) Matrix Spike Rec. outside of QC limits  
(anr) Analyte not requested

## MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D40798  
 Account: XTOKRWR - XTO Energy  
 Project: PCU 296-6A

QC Batch ID: MP8869  
 Matrix Type: SOLID

Methods: SW846 6010C  
 Units: mg/kg

Prep Date:

11/13/12

Metal	D40778-1 Original	MSD	Spikelot ICPALL2	% Rec	MSD RPD	QC Limit
Aluminum						
Antimony						
Arsenic						
Barium	3150	2830	219	-145.9(a	17.7	20
Beryllium						
Boron	anr					
Cadmium	0.0	46.3	54.8	84.5	9.9	20
Calcium						
Chromium	65.8	108	54.8	77.0	8.0	20
Cobalt						
Copper	14.3	62.1	54.8	87.2	9.2	20
Iron						
Lead	10.3	100	110	81.8	9.5	20
Lithium						
Magnesium						
Manganese						
Molybdenum						
Nickel	23.9	65.4	54.8	75.7	8.9	20
Phosphorus						
Potassium						
Selenium	0.64	94.8	110	85.9	9.3	20
Silicon						
Silver	0.13	19.4	21.9	87.9	9.8	20
Sodium						
Strontium						
Thallium						
Tin						
Titanium						
Uranium						
Vanadium						
Zinc	52.2	96.2	54.8	80.3	7.8	20

Associated samples MP8869: D40798-1

Results < IDL are shown as zero for calculation purposes  
 (\*) Outside of QC limits

MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D40798  
Account: XTOKRWR - XTO Energy  
Project: PCU 296-6A

QC Batch ID: MP8869  
Matrix Type: SOLID

Methods: SW846 6010C  
Units: mg/kg

Prep Date:

Metal

- (N) Matrix Spike Rec. outside of QC limits
- (anr) Analyte not requested
- (a) Spike amount low relative to the sample amount. Refer to lab control or spike blank for recovery information.

## SPIKE BLANK AND LAB CONTROL SAMPLE SUMMARY

Login Number: D40798  
 Account: XTOKRWR - XTO Energy  
 Project: PCU 296-6A

QC Batch ID: MP8869  
 Matrix Type: SOLID

Methods: SW846 6010C  
 Units: mg/kg

Prep Date:

11/13/12

Metal	BSP Result	Spikelot ICPALL2	% Rec	QC Limits
Aluminum				
Antimony				
Arsenic				
Barium	194	200	97.0	80-120
Beryllium				
Boron	anr			
Cadmium	45.0	50	90.0	80-120
Calcium				
Chromium	48.3	50	96.6	80-120
Cobalt				
Copper	45.7	50	91.4	80-120
Iron				
Lead	94.8	100	94.8	80-120
Lithium				
Magnesium				
Manganese				
Molybdenum				
Nickel	46.3	50	92.6	80-120
Phosphorus				
Potassium				
Selenium	93.5	100	93.5	80-120
Silicon				
Silver	18.9	20	94.5	80-120
Sodium				
Strontium				
Thallium				
Tin				
Titanium				
Uranium				
Vanadium				
Zinc	48.8	50	97.6	80-120

Associated samples MP8869: D40798-1

Results < IDL are shown as zero for calculation purposes  
 (\*) Outside of QC limits

SPIKE BLANK AND LAB CONTROL SAMPLE SUMMARY

Login Number: D40798  
Account: XTOKRWR - XTO Energy  
Project: PCU 296-6A

QC Batch ID: MP8869  
Matrix Type: SOLID

Methods: SW846 6010C  
Units: mg/kg

Prep Date:

Metal

(anr) Analyte not requested

## SERIAL DILUTION RESULTS SUMMARY

Login Number: D40798  
 Account: XTOKRWR - XTO Energy  
 Project: PCU 296-6A

QC Batch ID: MP8869  
 Matrix Type: SOLID

Methods: SW846 6010C  
 Units: ug/l

Prep Date:

11/13/12

Metal	D40778-1 Original	SDL 1:5	%DIF	QC Limits
Aluminum				
Antimony				
Arsenic				
Barium	27300	32600	19.4*(a)	0-10
Beryllium				
Boron	anr			
Cadmium	0.00	0.00	NC	0-10
Calcium				
Chromium	571	642	12.4*(a)	0-10
Cobalt				
Copper	124	113	8.9	0-10
Iron				
Lead	89.3	89.5	0.2	0-10
Lithium				
Magnesium				
Manganese				
Molybdenum				
Nickel	207	240	15.8*(a)	0-10
Phosphorus				
Potassium				
Selenium	5.60	0.00	100.0(b)	0-10
Silicon				
Silver	1.10	0.00	100.0(b)	0-10
Sodium				
Strontium				
Thallium				
Tin				
Titanium				
Uranium				
Vanadium				
Zinc	453	543	19.8*(a)	0-10

Associated samples MP8869: D40798-1

Results < IDL are shown as zero for calculation purposes  
 (\*) Outside of QC limits

SERIAL DILUTION RESULTS SUMMARY

Login Number: D40798  
Account: XTOKRWR - XTO Energy  
Project: PCU 296-6A

QC Batch ID: MP8869  
Matrix Type: SOLID

Methods: SW846 6010C  
Units: ug/l

Prep Date:

Metal

- (anr) Analyte not requested  
(a) Serial dilution indicates possible matrix interference.  
(b) Percent difference acceptable due to low initial sample concentration (< 50 times IDL).

BLANK RESULTS SUMMARY  
Part 2 - Method Blanks

Login Number: D40798  
Account: XTOKWR - XTO Energy  
Project: PCU 296-6A

QC Batch ID: MP8870  
Matrix Type: SOLID

Methods: SW846 6020A  
Units: mg/kg

Prep Date:

11/12/12

Metal	RL	IDL	MDL	MB raw	final
Aluminum	25	.22	.31		
Antimony	0.20	.0018	.0075		
Arsenic	0.10	.006	.06	0.0025	<0.10
Barium	1.0	.0065	.037		
Beryllium	0.10	.016	.09		
Boron	20	1.2	1.2		
Cadmium	0.050	.014	.021		
Calcium	200	7.9	8		
Chromium	1.0	.033	.19		
Cobalt	0.10	.0012	.015		
Copper	1.0	.017	.065		
Iron	20	.8	5		
Lead	0.25	.0011	.024		
Magnesium	50	.44	.85		
Manganese	0.50	.0043	.02		
Molybdenum	0.50	.018	.018		
Nickel	1.0	.0049	.011		
Phosphorus	30	1.4	3.6		
Potassium	100	9.8	10		
Selenium	0.20	.029	.14		
Silver	0.050	.0009	.0065		
Sodium	250	1.5	2.3		
Strontium	10	.036	.036		
Thallium	0.10	.00095	.0095		
Tin	5.0	.023	.34		
Titanium	1.0	.044	.1		
Uranium	0.25	.00085	.001		
Vanadium	2.0	.12	.21		
Zinc	5.0	.033	.35		

Associated samples MP8870: D40798-1

Results < IDL are shown as zero for calculation purposes  
(\*) Outside of QC limits  
(anr) Analyte not requested

## MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D40798  
 Account: XTOKRWR - XTO Energy  
 Project: PCU 296-6A

QC Batch ID: MP8870  
 Matrix Type: SOLID

Methods: SW846 6020A  
 Units: mg/kg

Prep Date:

11/12/12

Metal	D40778-1 Original MS	Spikelot ICPALL2	% Rec	QC Limits
Aluminum				
Antimony				
Arsenic	10.4	116	116	90.8    75-125
Barium				
Beryllium				
Boron				
Cadmium				
Calcium				
Chromium				
Cobalt				
Copper				
Iron				
Lead				
Magnesium				
Manganese				
Molybdenum				
Nickel				
Phosphorus				
Potassium				
Selenium				
Silver				
Sodium				
Strontium				
Thallium				
Tin				
Titanium				
Uranium				
Vanadium				
Zinc				

Associated samples MP8870: D40798-1

Results < IDL are shown as zero for calculation purposes  
 (\*) Outside of QC limits  
 (N) Matrix Spike Rec. outside of QC limits  
 (anr) Analyte not requested

## MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D40798  
 Account: XTOKRWR - XTO Energy  
 Project: PCU 296-6A

QC Batch ID: MP8870  
 Matrix Type: SOLID

Methods: SW846 6020A  
 Units: mg/kg

Prep Date:

11/12/12

Metal	D40778-1 Original	MSD	Spikelot ICPALL2	% Rec	MSD RPD	QC Limit
Aluminum						
Antimony						
Arsenic	10.4	104	110	85.4	10.9	20
Barium						
Beryllium						
Boron						
Cadmium						
Calcium						
Chromium						
Cobalt						
Copper						
Iron						
Lead						
Magnesium						
Manganese						
Molybdenum						
Nickel						
Phosphorus						
Potassium						
Selenium						
Silver						
Sodium						
Strontium						
Thallium						
Tin						
Titanium						
Uranium						
Vanadium						
Zinc						

Associated samples MP8870: D40798-1

Results < IDL are shown as zero for calculation purposes  
 (\*) Outside of QC limits  
 (N) Matrix Spike Rec. outside of QC limits  
 (anr) Analyte not requested

## SPIKE BLANK AND LAB CONTROL SAMPLE SUMMARY

Login Number: D40798  
 Account: XTOKRWR - XTO Energy  
 Project: PCU 296-6A

QC Batch ID: MP8870  
 Matrix Type: SOLID

Methods: SW846 6020A  
 Units: mg/kg

Prep Date: 11/12/12

Metal	BSP Result	Spikelot ICPALL2	% Rec	QC Limits
Aluminum				
Antimony				
Arsenic	92.2	100	92.2	80-120
Barium				
Beryllium				
Boron				
Cadmium				
Calcium				
Chromium				
Cobalt				
Copper				
Iron				
Lead				
Magnesium				
Manganese				
Molybdenum				
Nickel				
Phosphorus				
Potassium				
Selenium				
Silver				
Sodium				
Strontium				
Thallium				
Tin				
Titanium				
Uranium				
Vanadium				
Zinc				

Associated samples MP8870: D40798-1

Results < IDL are shown as zero for calculation purposes  
 (\*) Outside of QC limits  
 (anr) Analyte not requested

14.2.3  
**14**

## SERIAL DILUTION RESULTS SUMMARY

Login Number: D40798  
 Account: XTOKRWR - XTO Energy  
 Project: PCU 296-6A

QC Batch ID: MP8870  
 Matrix Type: SOLID

Methods: SW846 6020A  
 Units: ug/l

Prep Date:

11/12/12

Metal	D40778-1 Original	SDL 5:25	%DIF	QC Limits
Aluminum				
Antimony				
Arsenic	90.1	85.4	5.2	0-10
Barium				
Beryllium				
Boron				
Cadmium				
Calcium				
Chromium				
Cobalt				
Copper				
Iron				
Lead				
Magnesium				
Manganese				
Molybdenum				
Nickel				
Phosphorus				
Potassium				
Selenium				
Silver				
Sodium				
Strontium				
Thallium				
Tin				
Titanium				
Uranium				
Vanadium				
Zinc				

Associated samples MP8870: D40798-1

Results < IDL are shown as zero for calculation purposes  
 (\*) Outside of QC limits  
 (anr) Analyte not requested

14.2.4  
14

BLANK RESULTS SUMMARY  
Part 2 - Method Blanks

Login Number: D40798  
Account: XTOKRWR - XTO Energy  
Project: PCU 296-6A

QC Batch ID: MP8871  
Matrix Type: SOLID

Methods: SW846 7471B  
Units: mg/kg

Prep Date:

11/13/12

Metal	RL	IDL	MDL	MB raw	final
Mercury	0.083	.00088	.00075	-0.00024	<0.083

Associated samples MP8871: D40798-1

Results < IDL are shown as zero for calculation purposes  
(\*) Outside of QC limits  
(anr) Analyte not requested

MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D40798  
Account: XTOKWR - XTO Energy  
Project: PCU 296-6A

QC Batch ID: MP8871  
Matrix Type: SOLID

Methods: SW846 7471B  
Units: mg/kg

Prep Date: 11/13/12

Metal	D40797-1 Original MS	Spikelot HGWSR1	QC % Rec	QC Limits
Mercury	0.080	0.43	0.377	92.9 75-125

Associated samples MP8871: D40798-1

Results < IDL are shown as zero for calculation purposes

(\*) Outside of QC limits

(N) Matrix Spike Rec. outside of QC limits

(anr) Analyte not requested

MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D40798  
Account: XTOKWR - XTO Energy  
Project: PCU 296-6A

QC Batch ID: MP8871  
Matrix Type: SOLID

Methods: SW846 7471B  
Units: mg/kg

Prep Date:

11/13/12

Metal	D40797-1 Original	MSD	Spikelot HGWSR1	MSD % Rec	RPD	QC Limit
Mercury	0.080	0.40	0.371	86.3	7.2	20

Associated samples MP8871: D40798-1

Results < IDL are shown as zero for calculation purposes

(\*) Outside of QC limits

(N) Matrix Spike Rec. outside of QC limits

(anr) Analyte not requested

SPIKE BLANK AND LAB CONTROL SAMPLE SUMMARY

Login Number: D40798  
Account: XTOKRWR - XTO Energy  
Project: PCU 296-6A

QC Batch ID: MP8871  
Matrix Type: SOLID

Methods: SW846 7471B  
Units: mg/kg

Prep Date: 11/13/12

Metal	BSP Result	Spikelot HGWSR1	QC % Rec	QC Limits
Mercury	0.33	0.333	99.0	80-120

Associated samples MP8871: D40798-1

Results < IDL are shown as zero for calculation purposes  
(\*) Outside of QC limits  
(anr) Analyte not requested

BLANK RESULTS SUMMARY  
Part 2 - Method Blanks

Login Number: D40798  
Account: XTOKRWR - XTO Energy  
Project: PCU 296-6A

QC Batch ID: MP8877  
Matrix Type: AQUEOUS

Methods: SW846 6010C, USDA HANDBOOK 60  
Units: ug/l

Prep Date:

11/13/12

Metal	RL	IDL	MDL	MB raw	final
Aluminum	500	48	130		
Antimony	150	8.5	18		
Arsenic	130	22	42		
Barium	50	.5	9		
Beryllium	50	6.5	16		
Boron	250	5	22		
Cadmium	50	3	3		
Calcium	2000	27	80	16.5	<2000
Chromium	50	1.5	2.8		
Cobalt	25	2	2.1		
Copper	50	6	15		
Iron	350	6	100		
Lead	250	9.5	15		
Lithium	10	2.5			
Magnesium	1000	33	110	-0.50	<1000
Manganese	25	6	6		
Molybdenum	50	11	11		
Nickel	150	2.5	2.9		
Phosphorus	500	70	300		
Potassium	5000	310	750		
Selenium	250	24	55		
Silicon	250	15			
Silver	150	2	4.9		
Sodium	2000	30	490	147	<2000
Strontium	25	.2	7.5		
Thallium	50	15	43		
Tin	250	60			
Titanium	50	.5			
Uranium	250	11	23		
Vanadium	50	1	2.4		
Zinc	150	2.5	12		

Associated samples MP8877: D40798-1A

Results < IDL are shown as zero for calculation purposes  
(\*) Outside of QC limits

BLANK RESULTS SUMMARY  
Part 2 - Method Blanks

Login Number: D40798  
Account: XTOKRWR - XTO Energy  
Project: PCU 296-6A

QC Batch ID: MP8877  
Matrix Type: AQUEOUS

Methods: SW846 6010C, USDA HANDBOOK 60  
Units: ug/l

Prep Date:

Metal

(anr) Analyte not requested

## MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D40798  
 Account: XTOKRWR - XTO Energy  
 Project: PCU 296-6A

QC Batch ID: MP8877  
 Matrix Type: AQUEOUS

Methods: SW846 6010C, USDA HANDBOOK 60  
 Units: ug/l

Prep Date:

11/13/12

Metal	D40797-1A Original MS	Spikelot ICPALL2	% Rec	QC Limits
Aluminum				
Antimony				
Arsenic				
Barium				
Beryllium				
Boron				
Cadmium				
Calcium	14500	149000	125000	107.6
Chromium				
Cobalt				
Copper				
Iron				
Lead				
Lithium				
Magnesium	6860	128000	125000	96.9
Manganese				
Molybdenum				
Nickel				
Phosphorus				
Potassium				
Selenium				
Silicon				
Silver				
Sodium	56800	179000	125000	97.8
Strontium				
Thallium				
Tin				
Titanium				
Uranium				
Vanadium				
Zinc				

Associated samples MP8877: D40798-1A

Results < IDL are shown as zero for calculation purposes  
 (\*) Outside of QC limits

MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D40798  
Account: XTOKRWR - XTO Energy  
Project: PCU 296-6A

QC Batch ID: MP8877  
Matrix Type: AQUEOUS

Methods: SW846 6010C, USDA HANDBOOK 60  
Units: ug/l

Prep Date:

Metal

(N) Matrix Spike Rec. outside of QC limits  
(anr) Analyte not requested

## MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D40798  
 Account: XTOKRWR - XTO Energy  
 Project: PCU 296-6A

QC Batch ID: MP8877  
 Matrix Type: AQUEOUS

Methods: SW846 6010C, USDA HANDBOOK 60  
 Units: ug/l

Prep Date: 11/13/12

Metal	D40797-1A Original MSD	Spikelot ICPALL2	MSD % Rec	MSD RPD	QC Limit
Aluminum					
Antimony					
Arsenic					
Barium					
Beryllium					
Boron					
Cadmium					
Calcium	14500	153000	125000	110.8	2.6
Chromium					
Cobalt					
Copper					
Iron					
Lead					
Lithium					
Magnesium	6860	130000	125000	98.5	1.6
Manganese					
Molybdenum					
Nickel					
Phosphorus					
Potassium					
Selenium					
Silicon					
Silver					
Sodium	56800	180000	125000	98.6	0.6
Strontium					
Thallium					
Tin					
Titanium					
Uranium					
Vanadium					
Zinc					

Associated samples MP8877: D40798-1A

Results < IDL are shown as zero for calculation purposes  
 (\*) Outside of QC limits

MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D40798  
Account: XTOKRWR - XTO Energy  
Project: PCU 296-6A

QC Batch ID: MP8877  
Matrix Type: AQUEOUS

Methods: SW846 6010C, USDA HANDBOOK 60  
Units: ug/l

Prep Date:

Metal

(N) Matrix Spike Rec. outside of QC limits  
(anr) Analyte not requested

## SPIKE BLANK AND LAB CONTROL SAMPLE SUMMARY

Login Number: D40798  
 Account: XTOKRWR - XTO Energy  
 Project: PCU 296-6A

QC Batch ID: MP8877  
 Matrix Type: AQUEOUS

Methods: SW846 6010C, USDA HANDBOOK 60  
 Units: ug/l

Prep Date: 11/13/12

Metal	BSP Result	Spikelot ICPALL2	% Rec	QC Limits
Aluminum				
Antimony				
Arsenic				
Barium				
Beryllium				
Boron				
Cadmium				
Calcium	137000	125000	109.6	80-120
Chromium				
Cobalt				
Copper				
Iron				
Lead				
Lithium				
Magnesium	122000	125000	97.6	80-120
Manganese				
Molybdenum				
Nickel				
Phosphorus				
Potassium				
Selenium				
Silicon				
Silver				
Sodium	123000	125000	98.4	80-120
Strontium				
Thallium				
Tin				
Titanium				
Uranium				
Vanadium				
Zinc				

Associated samples MP8877: D40798-1A

Results < IDL are shown as zero for calculation purposes  
 (\*) Outside of QC limits

SPIKE BLANK AND LAB CONTROL SAMPLE SUMMARY

Login Number: D40798  
Account: XTOKRWR - XTO Energy  
Project: PCU 296-6A

QC Batch ID: MP8877  
Matrix Type: AQUEOUS

Methods: SW846 6010C, USDA HANDBOOK 60  
Units: ug/l

Prep Date:

Metal

(anr) Analyte not requested

## SERIAL DILUTION RESULTS SUMMARY

Login Number: D40798  
 Account: XTOKRWR - XTO Energy  
 Project: PCU 296-6A

QC Batch ID: MP8877  
 Matrix Type: AQUEOUS

Methods: SW846 6010C, USDA HANDBOOK 60  
 Units: ug/l

Prep Date: 11/13/12

Metal	D40797-1A	Original	SDL 1:5	%DIF	QC Limits
Aluminum					
Antimony					
Arsenic					
Barium					
Beryllium					
Boron					
Cadmium					
Calcium	2910	2990		2.8	0-10
Chromium					
Cobalt					
Copper					
Iron					
Lead					
Lithium					
Magnesium	1370	1400		2.4	0-10
Manganese					
Molybdenum					
Nickel					
Phosphorus					
Potassium					
Selenium					
Silicon					
Silver					
Sodium	11400	11600		2.6	0-10
Strontium					
Thallium					
Tin					
Titanium					
Uranium					
Vanadium					
Zinc					

Associated samples MP8877: D40798-1A

Results < IDL are shown as zero for calculation purposes  
 (\*) Outside of QC limits

SERIAL DILUTION RESULTS SUMMARY

Login Number: D40798  
Account: XTOKRWR - XTO Energy  
Project: PCU 296-6A

QC Batch ID: MP8877  
Matrix Type: AQUEOUS

Methods: SW846 6010C, USDA HANDBOOK 60  
Units: ug/l

Prep Date:

Metal

(anr) Analyte not requested



## General Chemistry

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### QC Data Summaries

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Includes the following where applicable:

- Method Blank and Blank Spike Summaries
- Duplicate Summaries
- Matrix Spike Summaries

METHOD BLANK AND SPIKE RESULTS SUMMARY  
GENERAL CHEMISTRY

Login Number: D40798  
Account: XTOKWR - XTO Energy  
Project: PCU 296-6A

Analyte	Batch ID	RL	MB Result	Units	Spike Amount	BSP Result	BSP %Recov	QC Limits
Chromium, Hexavalent	GP8655/GN17636	1.0	0.0	mg/kg	176	161	91.5	80-120%
Specific Conductivity	GP8677/GN17663			umhos/cm	9992	9790	98.0	90-110%
pH	GN17678			su	8.00	7.97	99.6	99.3-100.7%

Associated Samples:

Batch GP8655: D40798-1

Batch GP8677: D40798-1

Batch GN17678: D40798-1

(\*) Outside of QC limits

DUPLICATE RESULTS SUMMARY  
GENERAL CHEMISTRY

Login Number: D40798  
Account: XTOKWR - XTO Energy  
Project: PCU 296-6A

Analyte	Batch ID	QC Sample	Units	Original Result	DUP Result	RPD	QC Limits
Chromium, Hexavalent Redox Potential Vs H2	GP8655/GN17636 GN17680	D40715-1 D40797-1	mg/kg mv	0.0 138	0.0 140	0.0 1.4	0-20% 0-20%

Associated Samples:  
Batch GP8655: D40798-1  
Batch GN17680: D40798-1  
(\*) Outside of QC limits

MATRIX SPIKE RESULTS SUMMARY  
GENERAL CHEMISTRY

Login Number: D40798  
Account: XTOKWR - XTO Energy  
Project: PCU 296-6A

Analyte	Batch ID	QC Sample	Units	Original Result	Spike Amount	MS Result	%Rec	QC Limits
Chromium, Hexavalent	GP8655/GN17636	D40715-1	mg/kg	0.0	40	33.0	82.0	75-125%

Associated Samples:

Batch GP8655: D40798-1

(\*) Outside of QC limits

(N) Matrix Spike Rec. outside of QC limits

MATRIX SPIKE DUPLICATE RESULTS SUMMARY  
GENERAL CHEMISTRY

Login Number: D40798  
Account: XTOKWR - XTO Energy  
Project: PCU 296-6A

Analyte	Batch ID	QC Sample	Units	Original Result	Spike Amount	MSD Result	RPD	QC Limit
Chromium, Hexavalent	GP8655/GN17636	D40715-1	mg/kg	0.0	40	33.7	2.2	20%

Associated Samples:

Batch GP8655: D40798-1

(\*) Outside of QC limits

(N) Matrix Spike Rec. outside of QC limits