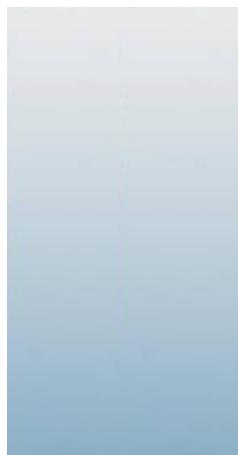




03/15/12



## Technical Report for

**XTO Energy**

**FRU 297-32A**

**1108-12A**

**Accutest Job Number: D32609**

**Sampling Date: 03/07/12**

### Report to:

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**Total number of pages in report: 134**



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, appearing to read "H. Madadian".

**Brad Madadian**  
**Laboratory Director**

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

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

This report shall not be reproduced, except in its entirety, without the written approval of Accutest Laboratories.

Test results relate only to samples analyzed.

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## Sample Summary

XTO Energy

**Job No:** D32609

FRU 297-32A

Project No: 1108-12A

Sample Number	Collected Date	Time By	Received	Matrix Code	Type	Client Sample ID
D32609-1	03/07/12	14:45 DS	03/12/12	SO	Soil	CUT 2 SUBLINER
D32609-1A	03/07/12	14:45 DS	03/12/12	SO	Soil	CUT 2 SUBLINER

---

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



## CASE NARRATIVE / CONFORMANCE SUMMARY

**Client:** XTO Energy

**Job No** D32609

**Site:** FRU 297-32A

**Report Date** 3/15/2012 4:32:42 PM

On 03/12/2012, 1 sample(s), 0 Trip Blank(s), and 0 Field Blank(s) were received at Accutest Mountain States (AMS) at a temperature of 3.6 °C. The samples were intact and properly preserved, unless noted below. An AMS Job Number of D32609 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> V5V1203
------------------	--------------------------

- All samples were analyzed within the recommended method holding time.
- Sample(s) D32550-1MS, D32550-1MSD were used as the QC samples indicated.
- All method blanks for this batch meet method specific criteria.
- The matrix spike duplicate (MSD) recovery(s) of Xylene (total) are outside control limits. Probable cause due to matrix interference.
- Sample(s) D32550-1MSD have surrogates outside control limits. Probable cause due to matrix interference.

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

<b>Matrix</b> SO	<b>Batch ID:</b> OP5528
------------------	-------------------------

- All samples were extracted and analyzed within the recommended method holding time.
- Sample(s) D32609-1MS, D32609-1MSD were used as the QC samples indicated.
- All method blanks for this batch meet method specific criteria.
- The RPD(s) for the MS and MSD recoveries of Acenaphthene, Chrysene, Dibenzo(a,h)anthracene, Fluorene, Indeno(1,2,3-cd)pyrene, Naphthalene are outside control limits for sample OP5528-MSD. Variability of recovery may be due to sample matrix/homogeneity.

### Volatiles by GC By Method SW846 8015B

<b>Matrix</b> SO	<b>Batch ID:</b> GGB858
------------------	-------------------------

- All samples were analyzed within the recommended method holding time.
- All method blanks for this batch meet method specific criteria.
- Sample(s) D32609-1MS, D32609-1MSD were used as the QC samples indicated.

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

<b>Matrix</b> SO	<b>Batch ID:</b> OP5519
------------------	-------------------------

- All samples were extracted D32554-1Awithin the recommended method holding time.
- All method blanks for this batch meet method specific criteria.
- Sample(s) D32595-1MSD, D32595-1MS, D32595-1MSD were used as the QC samples indicated.
- The matrix spike (MS) recovery(s) of TPH-DRO (C10-C28) are outside control limits. Outside control limits due to high level in sample relative to spike amount.
- Sample(s) OP5519-MS, OP5519-MSD have surrogates outside control limits. Probable cause due to matrix interference.
- OP5519-MS/MSD for o-Terphenyl: Outside control limits due to dilution.

## Metals By Method SW846 6010C

**Matrix** AQ

**Batch ID:** MP7061

- All samples were digested D32554-1A within the recommended method holding time.
- All method blanks for this batch meet method specific criteria.
- Sample(s) D32609-1AMS, D32609-1AMSD were used as the QC samples for the metals analysis.

**Matrix** SO

**Batch ID:** MP7055

- All samples were digested D32554-1A within the recommended method holding time.
- All method blanks for this batch meet method specific criteria.
- Sample(s) D32609-1MS, D32609-1MSD, D32609-1SDL were used as the QC samples for the metals analysis.
- The matrix spike (MS) recovery(s) of Barium are outside control limits. Spike amount low relative to the sample amount. Refer to lab control or spike blank for recovery information.
- The serial dilution RPD(s) for Cadmium, Selenium, Silver, Chromium, Lead, Nickel, Zinc are outside control limits for sample MP7055-SD1. Percent difference acceptable due to low initial sample concentration (< 50 times IDL).
- MP7055-SD1 for Chromium, Lead, Nickel, and Zinc: Serial dilution indicates possible matrix interference.

## Metals By Method SW846 6020A

**Matrix** SO

**Batch ID:** MP7056

- All samples were digested D32554-1A within the recommended method holding time.
- All method blanks for this batch meet method specific criteria.
- Sample(s) D32609-1MS, D32609-1MSD, D32609-1SDL were used as the QC samples for the metals analysis.
- The serial dilution RPD(s) for Arsenic are outside control limits for sample MP7056-SD1. Probable cause due to sample homogeneity.
- MP7056-SD1 for Arsenic: Serial dilution indicates possible matrix interference.

## Metals By Method SW846 7471B

**Matrix** SO

**Batch ID:** MP7050

- All samples were digested D32554-1A within the recommended method holding time.
- All method blanks for this batch meet method specific criteria.
- Sample(s) D32609-1MS, D32609-1MSD were used as the QC samples for the metals analysis.

## Wet Chemistry By Method ASTM D1498-76M

**Matrix** SO

**Batch ID:** GN14069

- Sample(s) D32609-1DUP 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:** GN14053

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

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

**Matrix** SO

**Batch ID:** R12120

- The data for SW846 3060/7196A M meets quality control requirements.
- D32609-1 for Chromium, Trivalent: Calculated as: (Chromium) - (Chromium, Hexavalent)

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

**Matrix** SO

**Batch ID:** GP6713

- All samples were prepared D32554-1A within the recommended method holding time.
- All method blanks for this batch meet method specific criteria.
- Sample(s) D32493-1DUP, D32493-1MS, D32493-1MSD were used as the QC samples for the Chromium, Hexavalent analysis.
- The matrix spike duplicate (MSD) recovery(s) of Chromium, Hexavalent are outside control limits. Probable cause due to matrix interference.
- GP6713-S2 for Chromium, Hexavalent: High RPD due to possible sample nonhomogeneity.

## Wet Chemistry By Method SW846 9045C

**Matrix** SO

**Batch ID:** GN14060

- The following samples were run outside of holding time for method SW846 9045C: D32609-1

## Wet Chemistry By Method USDA HANDBOOK 60

**Matrix** SO

**Batch ID:** MP7061

- D32609-1A 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.



## 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  
**Lab Sample ID:** D32609-1  
**Matrix:** SO - Soil  
**Method:** SW846 8260B  
**Project:** FRU 297-32A

**Date Sampled:** 03/07/12  
**Date Received:** 03/12/12  
**Percent Solids:** 94.4

	<b>File ID</b>	<b>DF</b>	<b>Analyzed</b>	<b>By</b>	<b>Prep Date</b>	<b>Prep Batch</b>	<b>Analytical Batch</b>
Run #1	5V19972.D	1	03/13/12	KV	n/a	n/a	V5V1203
Run #2							

	<b>Initial Weight</b>	<b>Final Volume</b>	<b>Methanol Aliquot</b>
Run #1	5.04 g	5.0 ml	100 ul
Run #2			

**Purgeable Aromatics**

<b>CAS No.</b>	<b>Compound</b>	<b>Result</b>	<b>RL</b>	<b>MDL</b>	<b>Units</b>	<b>Q</b>
71-43-2	Benzene	0.0469	0.056	0.024	mg/kg	J
108-88-3	Toluene	0.105	0.11	0.056	mg/kg	J
100-41-4	Ethylbenzene	ND	0.11	0.028	mg/kg	
1330-20-7	Xylene (total)	0.144	0.22	0.11	mg/kg	J

<b>CAS No.</b>	<b>Surrogate Recoveries</b>	<b>Run# 1</b>	<b>Run# 2</b>	<b>Limits</b>
2037-26-5	Toluene-D8	97%		61-130%
460-00-4	4-Bromofluorobenzene	104%		53-131%
17060-07-0	1,2-Dichloroethane-D4	107%		62-130%

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

Accutest Laboratories

**Report of Analysis**

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<b>Client Sample ID:</b>	CUT 2 SUBLINER	<b>Date Sampled:</b>	03/07/12
<b>Lab Sample ID:</b>	D32609-1	<b>Date Received:</b>	03/12/12
<b>Matrix:</b>	SO - Soil	<b>Percent Solids:</b>	94.4
<b>Method:</b>	SW846 8270C BY SIM	SW846 3546	
<b>Project:</b>	FRU 297-32A		
<b>File ID</b>	<b>DF</b>	<b>Analyzed</b>	<b>By</b>
Run #1	3G08526.D	1	03/14/12 DC
Run #2			
<b>Initial Weight</b>	<b>Final Volume</b>		
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	0.0088	0.0046	mg/kg	
120-12-7	Anthracene	ND	0.0088	0.0046	mg/kg	
56-55-3	Benzo(a)anthracene	ND	0.0088	0.0046	mg/kg	
50-32-8	Benzo(a)pyrene	ND	0.0088	0.0046	mg/kg	
205-99-2	Benzo(b)fluoranthene	ND	0.0088	0.0046	mg/kg	
207-08-9	Benzo(k)fluoranthene	ND	0.0088	0.0046	mg/kg	
218-01-9	Chrysene	ND	0.0088	0.0046	mg/kg	
53-70-3	Dibenz(a,h)anthracene	ND	0.0088	0.0046	mg/kg	
206-44-0	Fluoranthene	ND	0.0088	0.0046	mg/kg	
86-73-7	Fluorene	ND	0.0088	0.0046	mg/kg	
193-39-5	Indeno(1,2,3-cd)pyrene	ND	0.0088	0.0046	mg/kg	
91-20-3	Naphthalene	0.0327	0.012	0.011	mg/kg	
129-00-0	Pyrene	ND	0.0088	0.0046	mg/kg	
CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits		
4165-60-0	Nitrobenzene-d5	53%		10-145%		
321-60-8	2-Fluorobiphenyl	48%		10-130%		
1718-51-0	Terphenyl-d14	76%		22-130%		

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

Accutest Laboratories

**Report of Analysis**

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**Client Sample ID:** CUT 2 SUBLINER**Lab Sample ID:** D32609-1**Date Sampled:** 03/07/12**Matrix:** SO - Soil**Date Received:** 03/12/12**Method:** SW846 8015B**Percent Solids:** 94.4**Project:** FRU 297-32A

	<b>File ID</b>	<b>DF</b>	<b>Analyzed</b>	<b>By</b>	<b>Prep Date</b>	<b>Prep Batch</b>	<b>Analytical Batch</b>
Run #1	GB15297.D	1	03/13/12	SK	n/a	n/a	GGB858
Run #2							

	<b>Initial Weight</b>	<b>Final Volume</b>	<b>Methanol Aliquot</b>
Run #1	5.0 g	5.0 ml	100 ul
Run #2			

<b>CAS No.</b>	<b>Compound</b>	<b>Result</b>	<b>RL</b>	<b>MDL</b>	<b>Units</b>	<b>Q</b>
	TPH-GRO (C6-C10)	8.20	11	5.6	mg/kg	J
<b>CAS No.</b>	<b>Surrogate Recoveries</b>		<b>Run# 1</b>	<b>Run# 2</b>	<b>Limits</b>	
120-82-1	1,2,4-Trichlorobenzene	107%		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

Accutest Laboratories

**Report of Analysis**

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**Client Sample ID:** CUT 2 SUBLINER  
**Lab Sample ID:** D32609-1  
**Matrix:** SO - Soil  
**Method:** SW846-8015B SW846 3546  
**Project:** FRU 297-32A

**Date Sampled:** 03/07/12  
**Date Received:** 03/12/12  
**Percent Solids:** 94.4

	<b>File ID</b>	<b>DF</b>	<b>Analyzed</b>	<b>By</b>	<b>Prep Date</b>	<b>Prep Batch</b>	<b>Analytical Batch</b>
Run #1	FH002202.D	1	03/13/12	TR	03/12/12	OP5519	GFH113
Run #2							

	<b>Initial Weight</b>	<b>Final Volume</b>
Run #1	30.0 g	2.0 ml
Run #2		

<b>CAS No.</b>	<b>Compound</b>	<b>Result</b>	<b>RL</b>	<b>MDL</b>	<b>Units</b>	<b>Q</b>
	TPH-DRO (C10-C28)	115	14	9.2	mg/kg	
<b>CAS No.</b>	<b>Surrogate Recoveries</b>	<b>Run# 1</b>	<b>Run# 2</b>	<b>Limits</b>		
84-15-1	o-Terphenyl	65%		43-136%		

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**

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<b>Client Sample ID:</b>	CUT 2 SUBLINER	<b>Date Sampled:</b>	03/07/12
<b>Lab Sample ID:</b>	D32609-1	<b>Date Received:</b>	03/12/12
<b>Matrix:</b>	SO - Soil	<b>Percent Solids:</b>	94.4
<b>Project:</b>	FRU 297-32A		

**Metals Analysis**

Analyte	Result	RL	Units	DF	Prep	Analyzed By	Method	Prep Method
Arsenic	3.6	0.43	mg/kg	5	03/14/12	03/15/12 GJ	SW846 6020A <sup>3</sup>	SW846 3050B <sup>6</sup>
Barium	2100	1.1	mg/kg	1	03/14/12	03/14/12 JB	SW846 6010C <sup>2</sup>	SW846 3050B <sup>5</sup>
Cadmium	< 1.1	1.1	mg/kg	1	03/14/12	03/14/12 JB	SW846 6010C <sup>2</sup>	SW846 3050B <sup>5</sup>
Chromium	42.6	1.1	mg/kg	1	03/14/12	03/14/12 JB	SW846 6010C <sup>2</sup>	SW846 3050B <sup>5</sup>
Copper	13.5	1.1	mg/kg	1	03/14/12	03/14/12 JB	SW846 6010C <sup>2</sup>	SW846 3050B <sup>5</sup>
Lead	11.1	5.4	mg/kg	1	03/14/12	03/14/12 JB	SW846 6010C <sup>2</sup>	SW846 3050B <sup>5</sup>
Mercury	< 0.11	0.11	mg/kg	1	03/13/12	03/13/12 JM	SW846 7471B <sup>1</sup>	SW846 7471B <sup>4</sup>
Nickel	19.8	3.2	mg/kg	1	03/14/12	03/14/12 JB	SW846 6010C <sup>2</sup>	SW846 3050B <sup>5</sup>
Selenium	< 5.4	5.4	mg/kg	1	03/14/12	03/14/12 JB	SW846 6010C <sup>2</sup>	SW846 3050B <sup>5</sup>
Silver	< 3.2	3.2	mg/kg	1	03/14/12	03/14/12 JB	SW846 6010C <sup>2</sup>	SW846 3050B <sup>5</sup>
Zinc	43.1	3.2	mg/kg	1	03/14/12	03/14/12 JB	SW846 6010C <sup>2</sup>	SW846 3050B <sup>5</sup>

- (1) Instrument QC Batch: MA2256
- (2) Instrument QC Batch: MA2258
- (3) Instrument QC Batch: MA2262
- (4) Prep QC Batch: MP7050
- (5) Prep QC Batch: MP7055
- (6) Prep QC Batch: MP7056

RL = Reporting Limit

**Report of Analysis**

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**Client Sample ID:** CUT 2 SUBLINER**Lab Sample ID:** D32609-1**Matrix:** SO - Soil**Date Sampled:** 03/07/12**Date Received:** 03/12/12**Percent Solids:** 94.4**Project:** FRU 297-32A**General Chemistry**

Analyte	Result	RL	Units	DF	Analyzed	By	Method
Chromium, Hexavalent	< 1.0	1.0	mg/kg	1	03/14/12	CJ	SW846 3060A/7196A
Chromium, Trivalent <sup>a</sup>	42.6	2.1	mg/kg	1	03/14/12 22:54	JB	SW846 3060/7196A M
Redox Potential Vs H2	391		mv	1	03/13/12	JD	ASTM D1498-76M
Solids, Percent	94.4		%	1	03/13/12	SWT	SM19 2540B M
Specific Conductivity	1340	1.0	umhos/cm	1	03/14/12	JD	DEPT.OF AG, BOOK N9
pH	9.89		su	1	03/13/12 10:30	CT	SW846 9045C

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

RL = Reporting Limit

**Report of Analysis**

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<b>Client Sample ID:</b>	CUT 2 SUBLINER	<b>Date Sampled:</b>	03/07/12
<b>Lab Sample ID:</b>	D32609-1A	<b>Date Received:</b>	03/12/12
<b>Matrix:</b>	SO - Soil	<b>Percent Solids:</b>	94.4
<b>Project:</b>	FRU 297-32A		

**SAR Metals Analysis**

Analyte	Result	RL	Units	DF	Prep	Analyzed By	Method	Prep Method
Calcium	18.0	2.0	mg/l	1	03/14/12	03/14/12 JB	SW846 6010C <sup>1</sup>	EPA 200.7 <sup>2</sup>
Magnesium	14.9	1.0	mg/l	1	03/14/12	03/14/12 JB	SW846 6010C <sup>1</sup>	EPA 200.7 <sup>2</sup>
Sodium	248	2.0	mg/l	1	03/14/12	03/14/12 JB	SW846 6010C <sup>1</sup>	EPA 200.7 <sup>2</sup>

(1) Instrument QC Batch: MA2258

(2) Prep QC Batch: MP7061

RL = Reporting Limit

**Report of Analysis**

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<b>Client Sample ID:</b>	CUT 2 SUBLINER	<b>Date Sampled:</b>	03/07/12
<b>Lab Sample ID:</b>	D32609-1A	<b>Date Received:</b>	03/12/12
<b>Matrix:</b>	SO - Soil	<b>Percent Solids:</b>	94.4
<b>Project:</b>	FRU 297-32A		

**General Chemistry**

Analyte	Result	RL	Units	DF	Analyzed	By	Method
Sodium Adsorption Ratio <sup>a</sup>	10.5		ratio	1	03/14/12 14:14	JB	USDA HANDBOOK 60

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

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RL = Reporting Limit



## Misc. Forms

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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

Accutest Laboratories Mountain States  
4036 Youngfield Street Wheat Ridge, Co 80033  
TEL. 303-425-6021 877-737-4521  
FAX 303-425-6021

FED-EX Tracking #	Bottle Order Control #
Accutest Quote #	Accutest Job # D32609

Client / Reporting Information		Project Information		Requested Analysis (see TEST CODE sheet)		Matrix Codes	
Company Name <b>KRW CONSULTING</b>	Project Name <b>XTO FRU 207-32A</b>	Street: <b>8000 W 14TH AVE STE 200</b>	Billing Information (If different from Report to) <b>XTO ENERGY</b>				DW - Drinking Water GW - Ground Water WW - Water SW - Surface Water SO - Soil SL - Sludge SED - Sediment OL - Oil LIQ - Other Liquid AIR - Air SOL - Other Solid WP - Wipe FB - Field Blank EB - Equipment Blank RB - Rinse Blank TB - Trip Blank
City State Zip <b>LAKWOOD CO 80214</b>	Project# <b>1109-12A</b>	City: <b>RIFLE CO 81650</b>	Street Address <b>21459 CRS</b>	City State Zip <b>RIFLE CO 81650</b>	PO# <b>JESSICA DOOLING</b>		
Project Contact Email <b>DMAYNE KNUDSON</b>	Phone# <b>970 488 1098</b>	Client PO#	Project Manager <b>JOE HESS</b>	Attention:			LAB USE ONLY <b>D1</b> <i>3/12/12</i>
Sampler(s) Name(s) <b>DAVID SANDERS</b>	Phone # <b>970 488 1098</b>						
Collection		Number of preserved Bottles					
Accutest Sample #	Field ID / Point of Collection <b>CUT 2 SUBLINER</b>	Date <b>3-7-12 14:45</b>	Time <b>DUS</b>	Sampled by <b>SD</b>	Matrix <b>HCl</b>	# of bottles <b>5</b>	HCl NaOH HNCO H2SO4 NONE DI Water MEOH ENCORE Bisulfite
							X TABLE 910
Date Deliverable Information						Comments / Special Instructions <b>PLEASE EMAIL RESULTS TO KRW PICEANCE TEAM</b>	
Turnaround Time (Business days)		Approved By (Accutest PM): / Date:		<input type="checkbox"/> Commercial "A" (Level 1) <input type="checkbox"/> Commercial "B" (Level 2) <input type="checkbox"/> Commercial "B" + Narrative <input checked="" type="checkbox"/> FULLT1 (Level 3+4)  <input type="checkbox"/> State Forms <input type="checkbox"/> EDD Format <input checked="" type="checkbox"/> PDF			
<input type="checkbox"/> Std. 10 Business Days <input checked="" type="checkbox"/> Std. 5 Business Days (By Contract only) <input type="checkbox"/> 5 Day R/H SH <input checked="" type="checkbox"/> 3 Day EMERGENC <input type="checkbox"/> 2 Day EMERGENC <input type="checkbox"/> 1 Day EMERGENC							
Emergency & Rush T/A data available VIA LabLink.							
Sample Custody must be documented below outlining samples change possession, including courier delivery.							
Relinquished by Supplier: <b>1</b>	Date Time: <b>3/8/12 1730</b>	Received By: <b>AMERICAN COURIER</b>	Relinquished By: <b>2</b>	Date Time: <b>3/8/12 1730</b>	Received By: <b>AMERICAN COURIER</b>	Received By: <b>2</b>	Date Time: <b>3/12/12 1330</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>	Received By: <b>4</b>	Date Time:
Relinquished by: <b>5</b>	Date Time:	Received By: <b>5</b>	Custody Seal #		Intact <input type="checkbox"/> Preserved where applicable <input type="checkbox"/> Not intact <input type="checkbox"/>	On Ice <b>✓</b>	Cooler Temp. <b>3.6</b>

This  
Date

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**D32609: Chain of Custody**  
**Page 1 of 2**



## Accutest Laboratories Sample Receipt Summary

Accutest Job Number: D32609

Client: KRW CONSULTING

Immediate Client Services Action Required: No

Date / Time Received: 3/12/2012 1:30:00 PM

No. Coolers:

1

Client Service Action Required at Login: No

Project: XTO FRU 297-32A

Airbill #'s: HD/CO

**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:    
 2. Cooler temp verification: Infared gun  
 3. Cooler media: Ice (bag)

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

1. Trip Blank present / cooler:    
 2. Trip Blank listed on COC:    
 3. Samples preserved properly:    
 4. VOCs headspace free:

**Sample Integrity - Documentation****Y or N**

1. Sample labels present on bottles:    
 2. Container labeling complete:    
 3. Sample container label / COC agree:

**Sample Integrity - Condition****Y or N**

1. Sample recvd within HT:    
 2. All containers accounted for:    
 3. Condition of sample: Intact

**Sample Integrity - Instructions****Y or N****N/A**

1. Analysis requested is clear:    
 2. Bottles received for unspecified tests:    
 3. Sufficient volume rec'd for analysis:    
 4. Compositing instructions clear:     
 5. Filtering instructions clear:

Comments

Accutest Laboratories  
V:(303) 425-60214036 Youngfield Street  
F: (303) 425-6854Wheat Ridge, CO  
www.accutest.com

4.1

4

**D32609: Chain of Custody****Page 2 of 2**



## GC/MS Volatiles

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### QC Data Summaries

Includes the following where applicable:

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

**Method Blank Summary**

**Job Number:** D32609  
**Account:** XTOKWR XTO Energy  
**Project:** FRU 297-32A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
V5V1203-MB	5V19961.D	1	03/13/12	KV	n/a	n/a	V5V1203

The QC reported here applies to the following samples:

**Method:** SW846 8260B

D32609-1

CAS No.	Compound	Result	RL	MDL	Units	Q
71-43-2	Benzene	ND	50	22	ug/kg	
100-41-4	Ethylbenzene	ND	100	25	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	104%
460-00-4	4-Bromofluorobenzene	99%
17060-07-0	1,2-Dichloroethane-D4	115%

## Blank Spike Summary

Page 1 of 1

Job Number: D32609

Account: XTOKWR XTO Energy

Project: FRU 297-32A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
V5V1203-BS	5V19962.D	1	03/13/12	KV	n/a	n/a	V5V1203

The QC reported here applies to the following samples:

Method: SW846 8260B

D32609-1

CAS No.	Compound	Spike ug/kg	BSP ug/kg	BSP %	Limits
71-43-2	Benzene	50	49.4	99	70-130
100-41-4	Ethylbenzene	50	49.1	98	70-130
108-88-3	Toluene	50	47.5	95	70-130
1330-20-7	Xylene (total)	150	144	96	70-130

CAS No.	Surrogate Recoveries	BSP	Limits
2037-26-5	Toluene-D8	93%	61-130%
460-00-4	4-Bromofluorobenzene	102%	53-131%
17060-07-0	1,2-Dichloroethane-D4	97%	62-130%

# Matrix Spike/Matrix Spike Duplicate Summary

Page 1 of 1

Job Number: D32609

Account: XTOKWR XTO Energy

Project: FRU 297-32A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
D32550-1MS	5V19964.D	1	03/13/12	KV	n/a	n/a	V5V1203
D32550-1MSD	5V19965.D	1	03/13/12	KV	n/a	n/a	V5V1203
D32550-1	5V19963.D	1	03/13/12	KV	n/a	n/a	V5V1203

The QC reported here applies to the following samples:

Method: SW846 8260B

D32609-1

CAS No.	Compound	D32550-1		Spike	MS	MS	MSD	MSD	Limits	
		ug/kg	Q	ug/kg	ug/kg	%	ug/kg	%	RPD	Rec/RPD
71-43-2	Benzene	384		5830	6590	106	7520	122	13	70-134/30
100-41-4	Ethylbenzene	1470		5830	7400	102	8300	117	11	70-137/30
108-88-3	Toluene	5170		5830	11000	100	12300	122	11	70-130/30
1330-20-7	Xylene (total)	28000		17500	46200	104	51400	134* a	11	61-131/30

CAS No.	Surrogate Recoveries	MS	MSD	D32550-1	Limits
2037-26-5	Toluene-D8	99%	109%	89%	61-130%
460-00-4	4-Bromofluorobenzene	126%	138% *	101%	53-131%
17060-07-0	1,2-Dichloroethane-D4	95%	105%	87%	62-130%

(a) Outside control limits due to possible matrix interference.



## GC/MS Volatiles

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Raw Data

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## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\V5031312.S\  
 Data File : 5V19972.D  
 Acq On : 13 Mar 2012 4:23 pm  
 Operator : KOROUSHV  
 Sample : D32609-1  
 Misc : MS3551,V5V1203,5.038,,100,5,1  
 ALS Vial : 14 Sample Multiplier: 1

Quant Time: Mar 14 10:06:32 2012  
 Quant Method : C:\msdchem\1\METHODS\V5AP1186TVH1186.M  
 Quant Title : 8260  
 QLast Update : Fri Mar 02 14:22:16 2012  
 Response via : Initial Calibration

6.1.1

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
2) Pentafluorobenzene	11.647	168	189646	50.00	ug/l	0.00
35) 1,4-Difluorobenzene	12.446	114	342727	50.00	ug/l	0.00
53) Chlorobenzene-d5	15.095	117	522117	50.00	ug/l	0.00
74) 1,4-Dichlorobenzene-d4	17.070	152	372229	50.00	ug/l	0.00

System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	12.035	102	43611	53.42	ug/l	0.00
Spiked Amount	50.000	Range	70 - 130	Recovery	=	106.84%
61) Toluene-d8	13.850	98	816665	48.38	ug/l	0.00
Spiked Amount	50.000	Range	70 - 130	Recovery	=	96.76%
69) 4-Bromofluorobenzene	16.043	95	363474	52.13	ug/l	0.00
Spiked Amount	50.000	Range	70 - 130	Recovery	=	104.26%

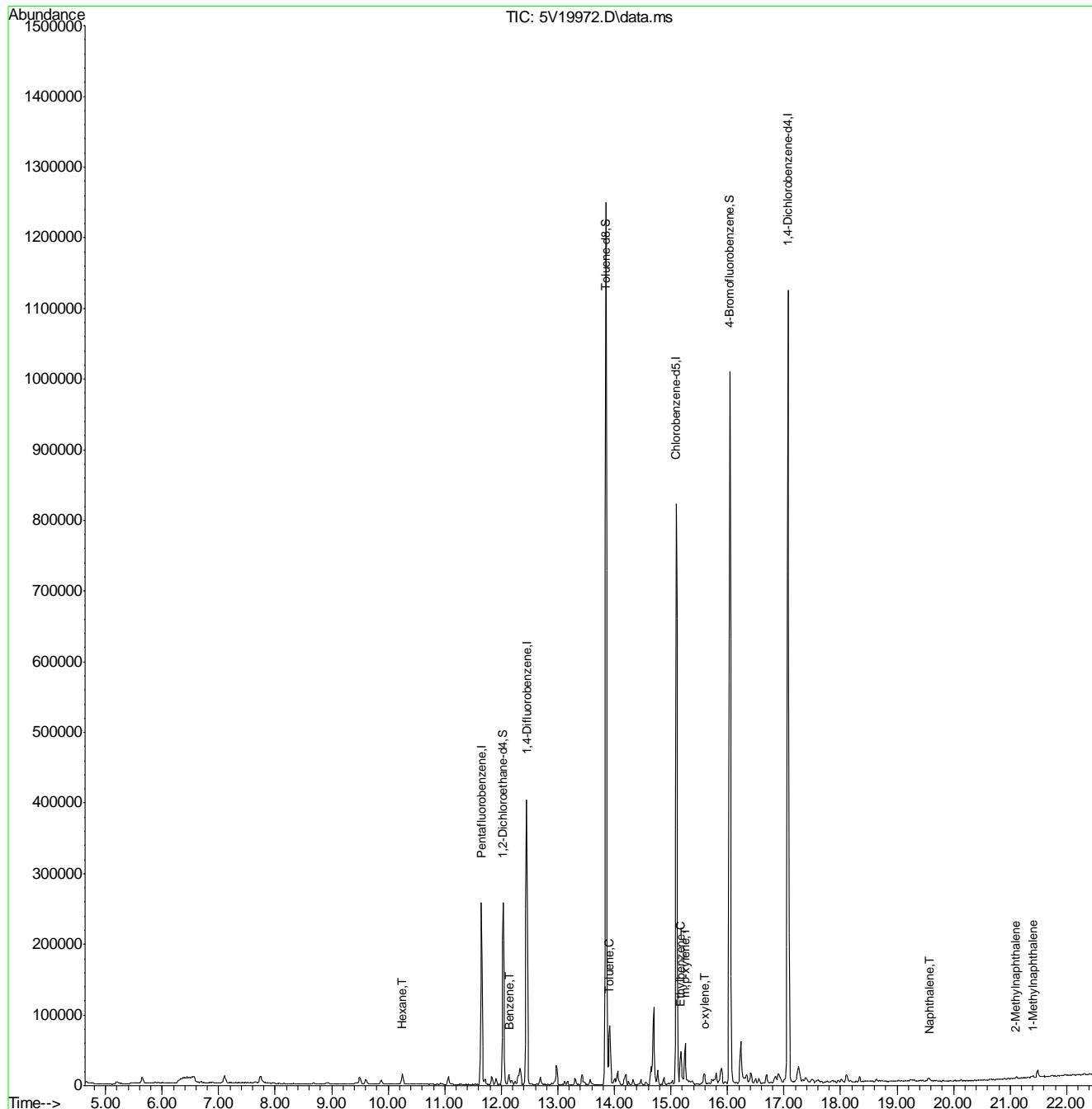
Target Compounds					QValue
41) Hexane	10.243	57	8080	2.29	ug/l 100
50) Benzene	12.138	78	14222	0.84	ug/l 100
62) Toluene	13.908	92	24261	1.89	ug/l 96
66) Ethylbenzene	15.175	91	9917	0.41	ug/l 100
72) m,p-xylene	15.255	106	18700	1.91	ug/l 98
73) o-xylene	15.597	106	2176	0.68	ug/l 84
91) Naphthalene	19.570	128	4594	1.49	ug/l 100
94) 2-Methylnaphthalene	21.100	142	1706	2.15	ug/l 97
95) 1-Methylnaphthalene	21.408	142	1092	1.99	ug/l # 85

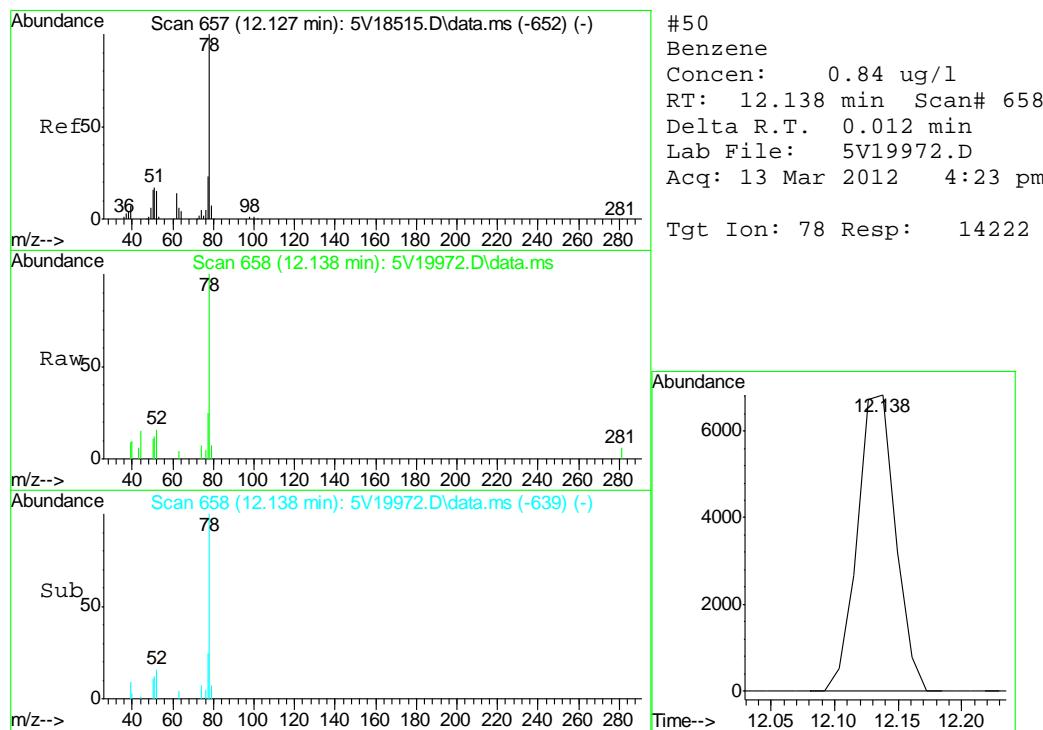
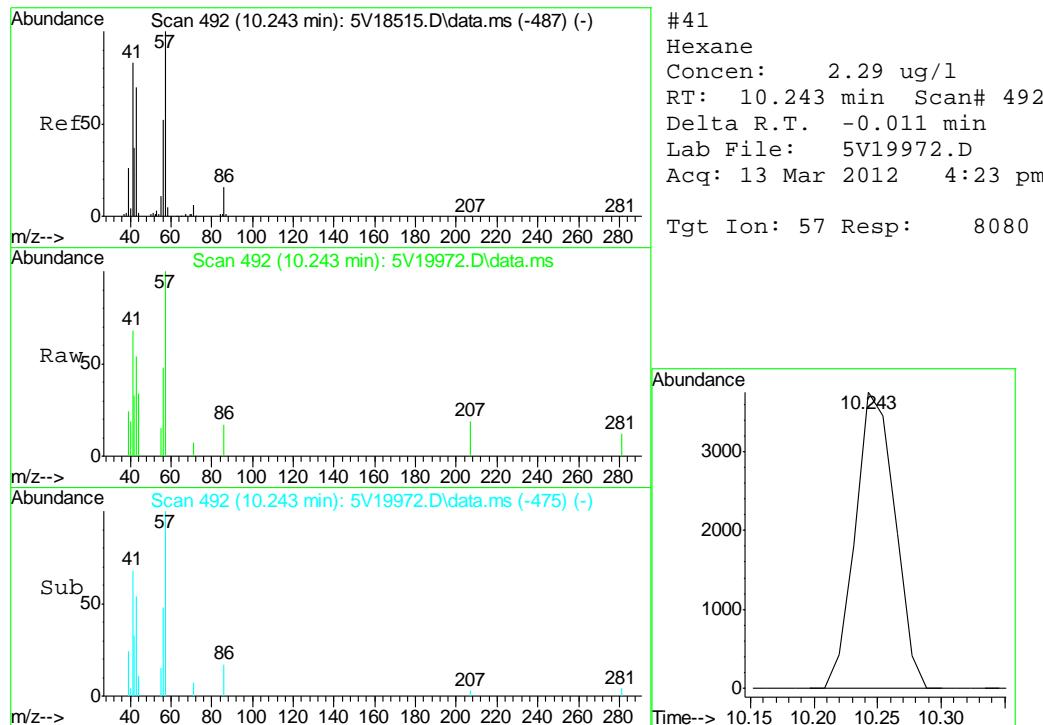
(#) = qualifier out of range (m) = manual integration (+) = signals summed

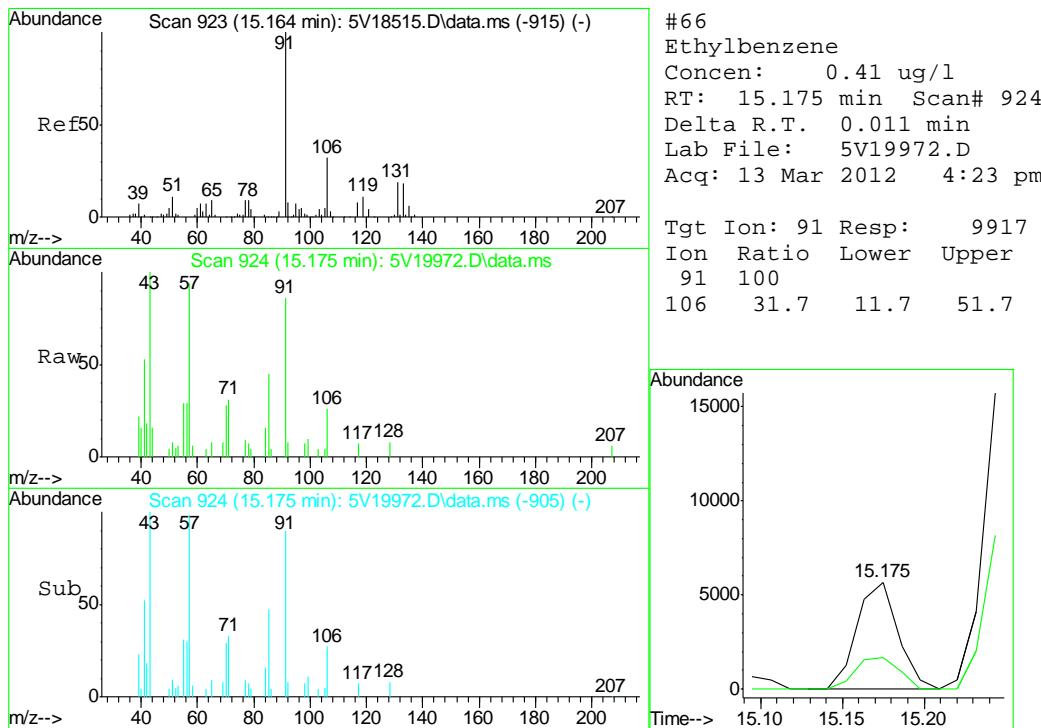
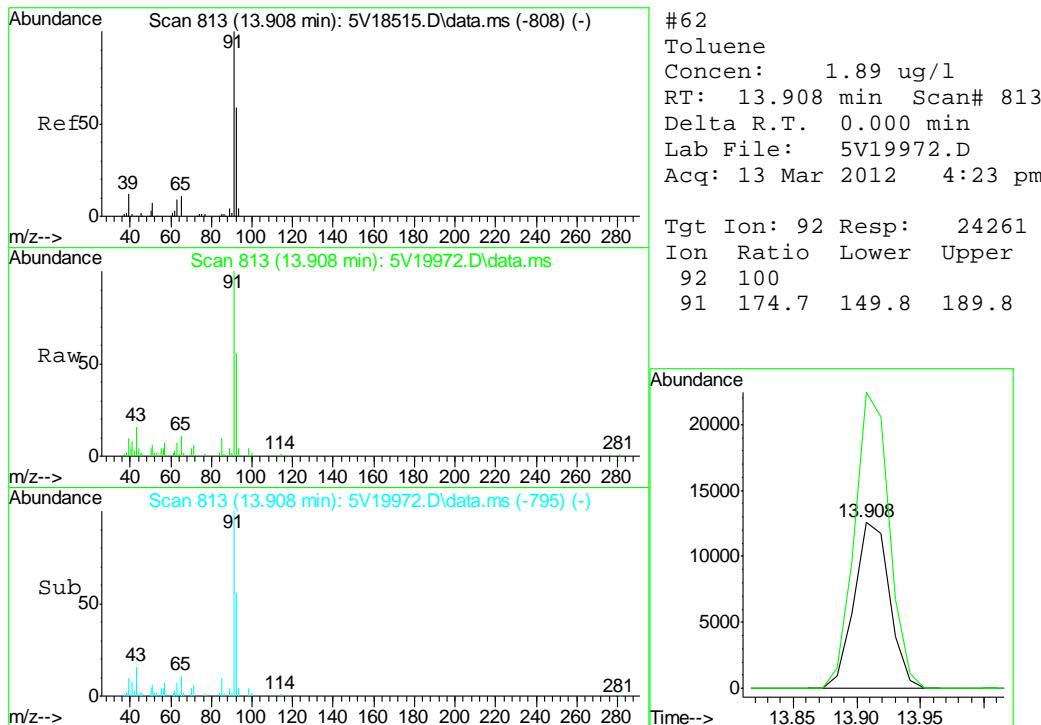
## Quantitation Report (QT Reviewed)

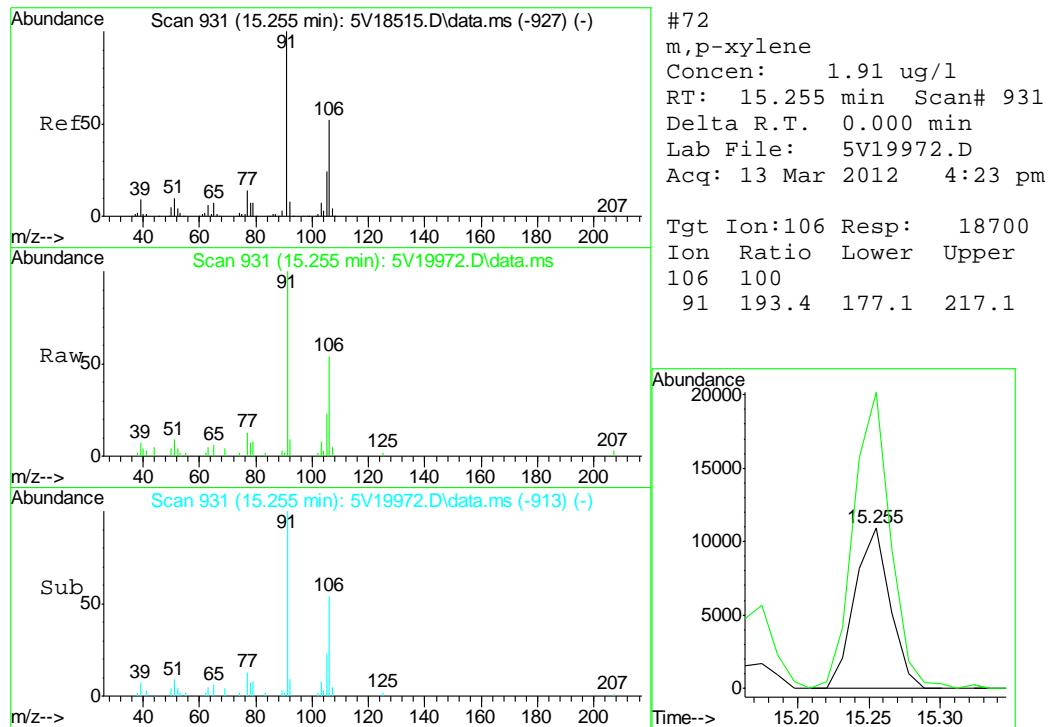
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 Data File : 5V19972.D  
 Acq On : 13 Mar 2012 4:23 pm  
 Operator : KOROUSHV  
 Sample : D32609-1  
 Misc : MS3551,V5V1203,5.038,,100,5,1  
 ALS Vial : 14 Sample Multiplier: 1

Quant Time: Mar 14 10:06:32 2012  
 Quant Method : C:\msdchem\1\METHODS\V5AP1186TVH1186.M  
 Quant Title : 8260  
 QLast Update : Fri Mar 02 14:22:16 2012  
 Response via : Initial Calibration



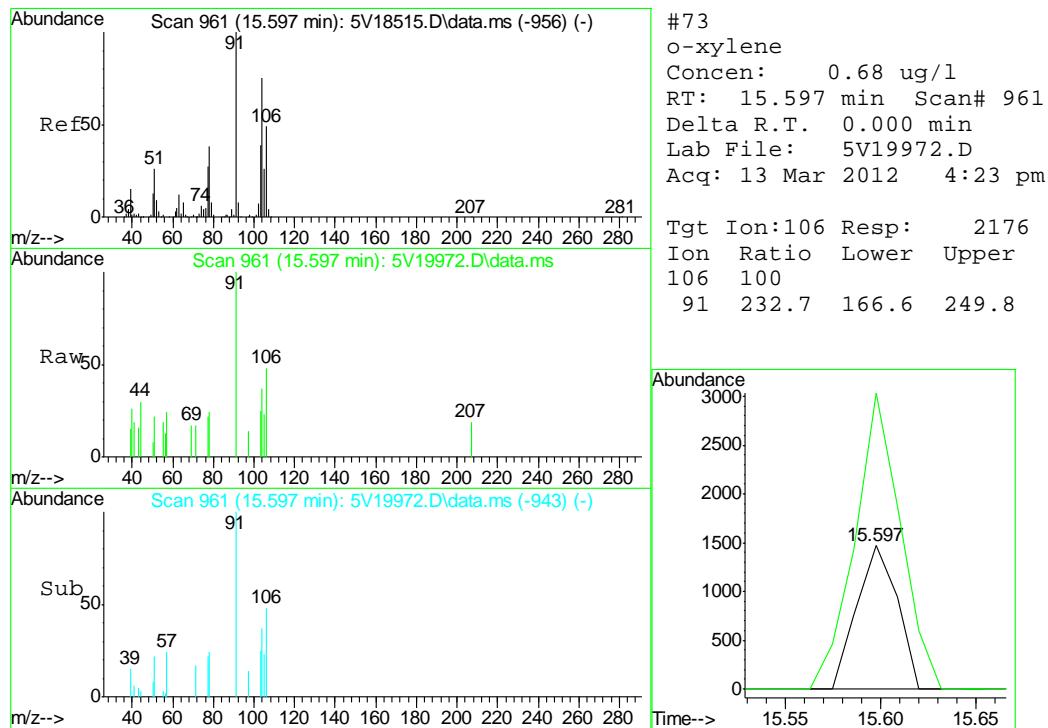


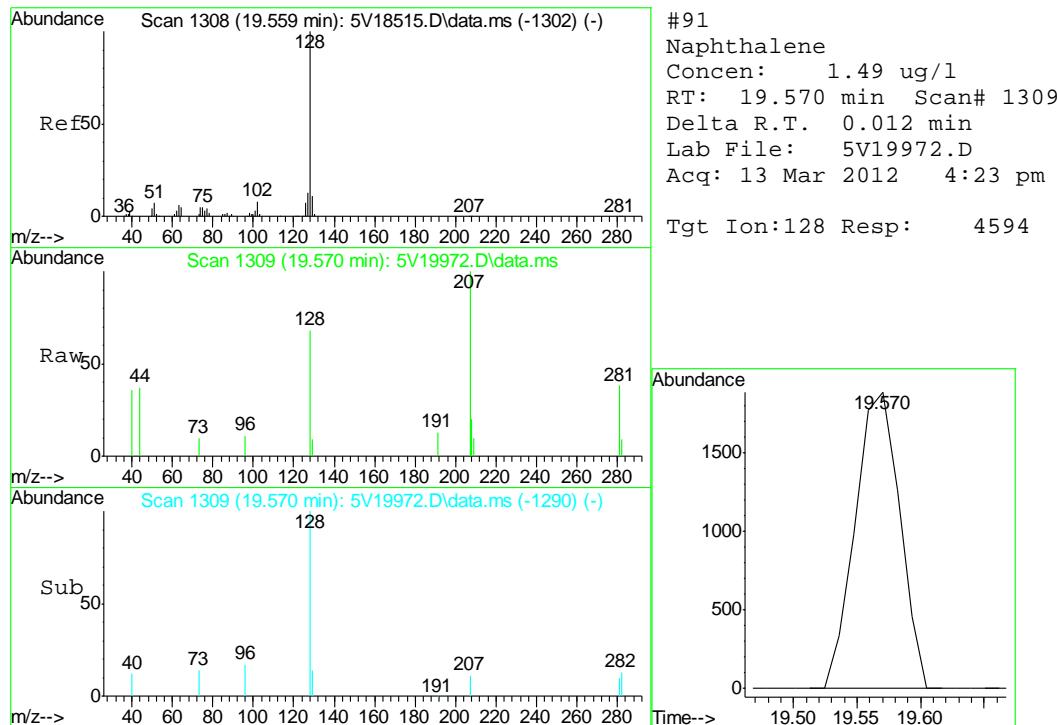




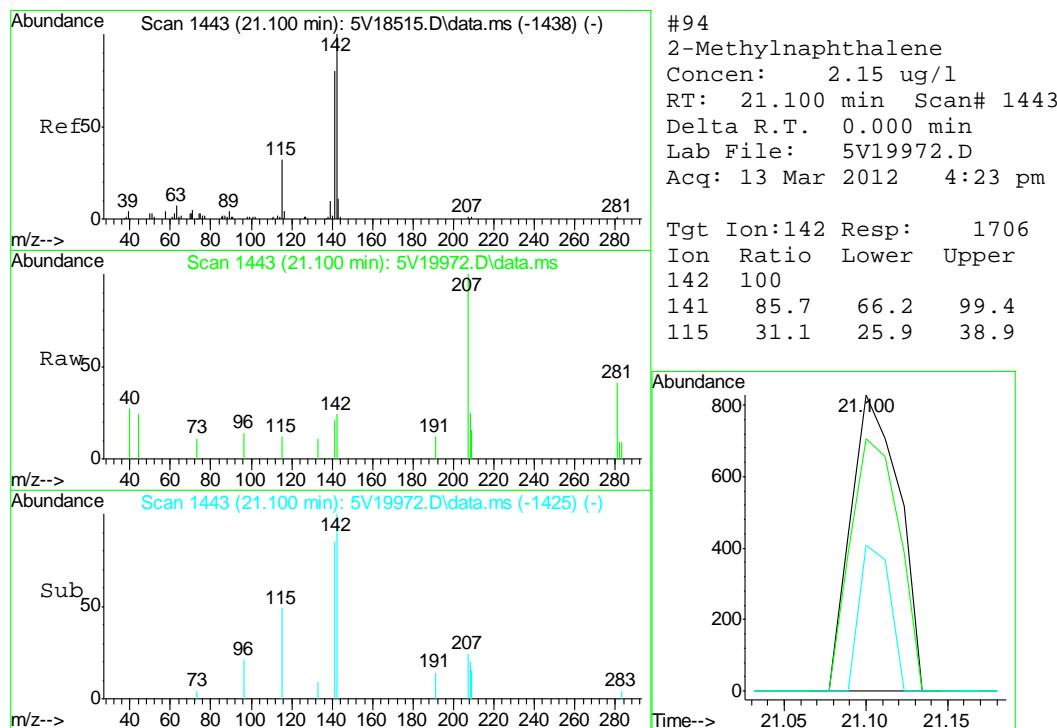
6.1.1

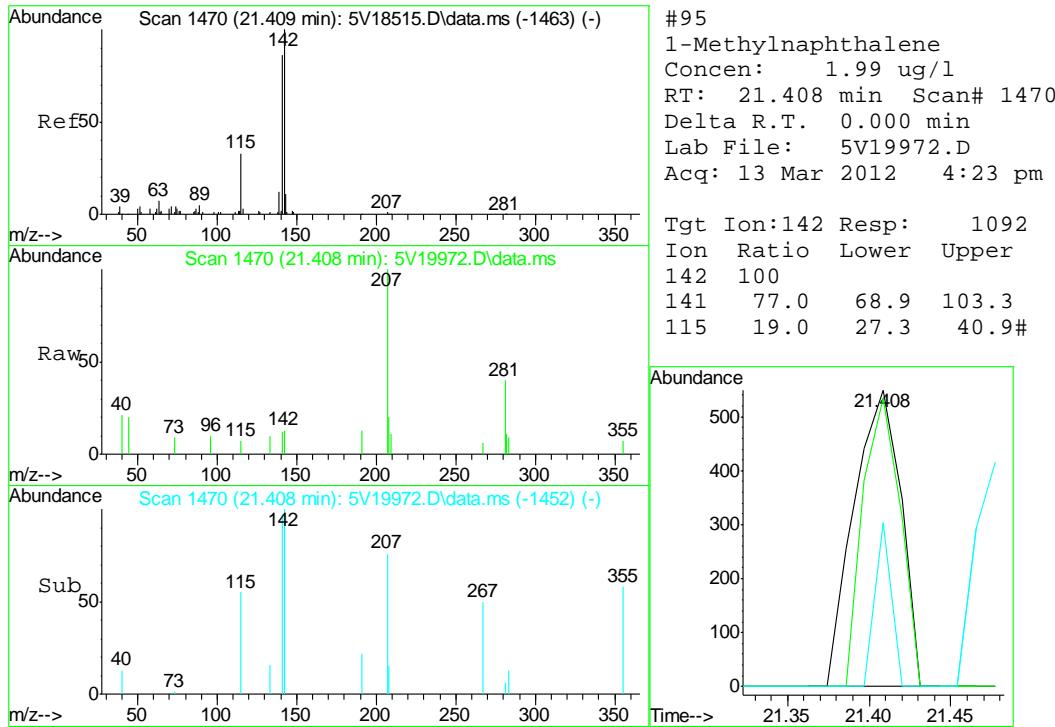
6





6.1.1





## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\V5031312.S\  
 Data File : 5V19961.D  
 Acq On : 13 Mar 2012 10:14 am  
 Operator : KOROUSHV  
 Sample : MB  
 Misc : MS3551,V5V1203,5.00,,100,5,1  
 ALS Vial : 3 Sample Multiplier: 1

Quant Time: Mar 14 09:51:19 2012  
 Quant Method : C:\msdchem\1\METHODS\V5AP1186TVH1186.M  
 Quant Title : 8260  
 QLast Update : Fri Mar 02 14:22:16 2012  
 Response via : Initial Calibration

6.2.1

6

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
2) Pentafluorobenzene	11.647	168	192012	50.00	ug/l	0.00
35) 1,4-Difluorobenzene	12.446	114	342802	50.00	ug/l	0.00
53) Chlorobenzene-d5	15.095	117	507384	50.00	ug/l	0.00
74) 1,4-Dichlorobenzene-d4	17.070	152	334611	50.00	ug/l	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
33) 1,2-Dichloroethane-d4	12.035	102	47123	57.25	ug/l	0.00
Spiked Amount 50.000	Range 70 - 130		Recovery	=	114.50%	
61) Toluene-d8	13.850	98	857074	52.25	ug/l	0.00
Spiked Amount 50.000	Range 70 - 130		Recovery	=	104.50%	
69) 4-Bromofluorobenzene	16.043	95	335704	49.54	ug/l	0.00
Spiked Amount 50.000	Range 70 - 130		Recovery	=	99.08%	

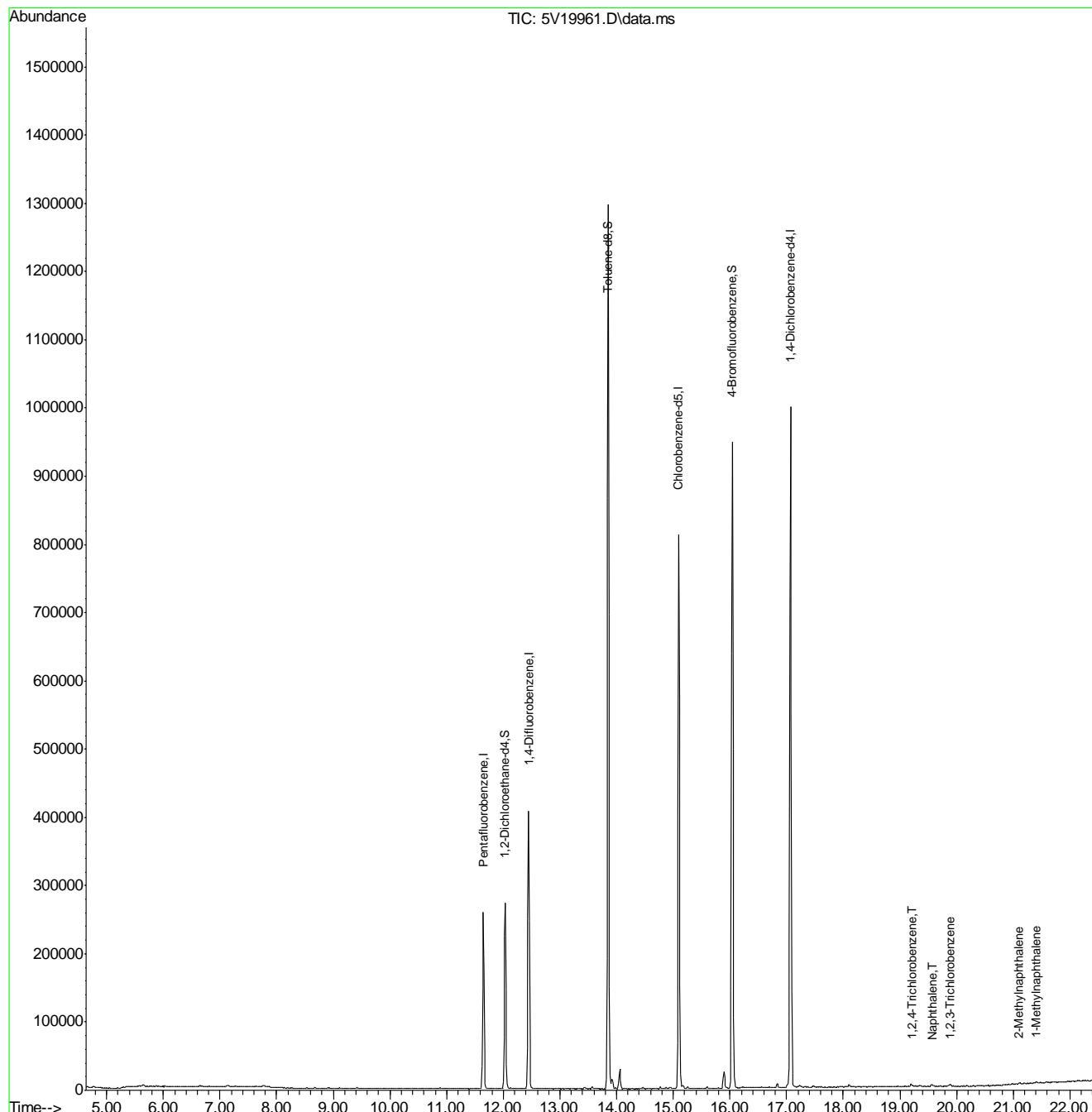
Target Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)	QValue
90) 1,2,4-Trichlorobenzene	19.205	180	1996	0.83	ug/l	#	80
91) Naphthalene	19.570	128	5625	1.59	ug/l		100
93) 1,2,3-Trichlorobenzene	19.879	180	2428	0.86	ug/l	#	89
94) 2-Methylnaphthalene	21.100	142	1211	2.05	ug/l	#	84
95) 1-Methylnaphthalene	21.397	142	1655	2.14	ug/l	#	85

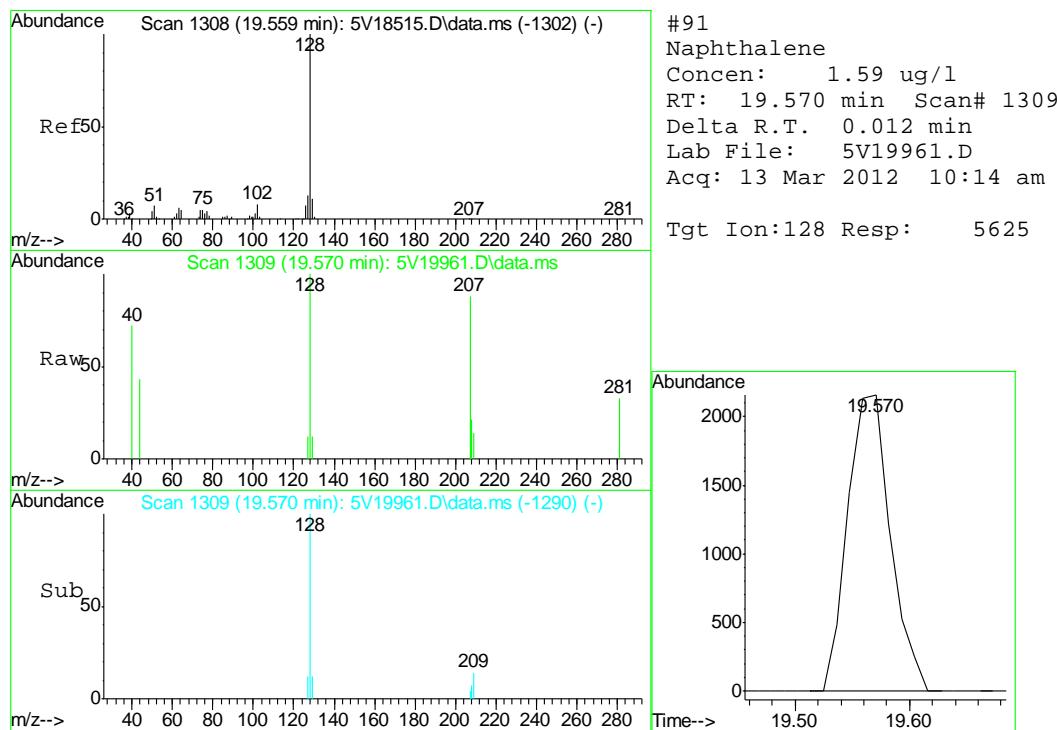
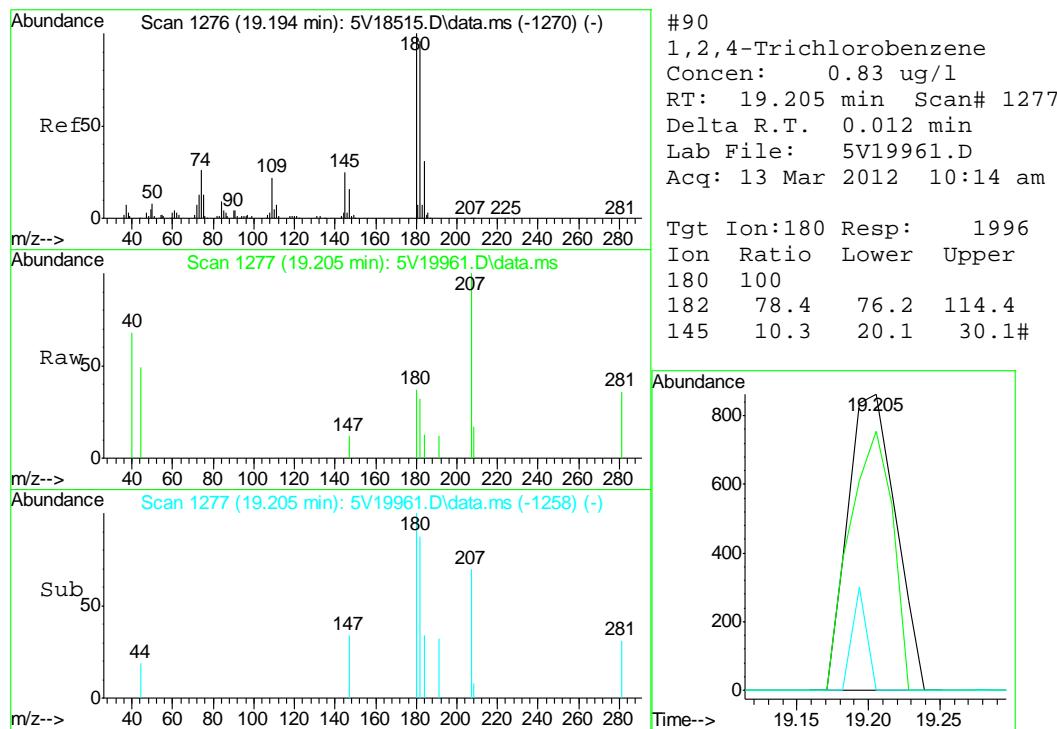
(#) = qualifier out of range (m) = manual integration (+) = signals summed

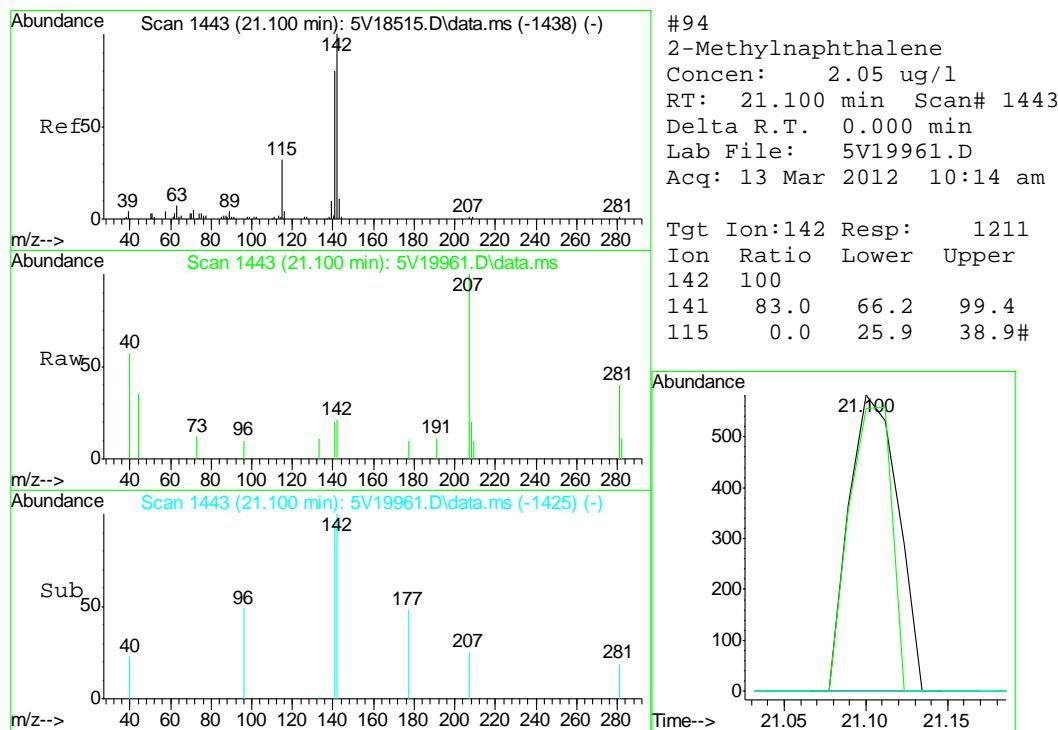
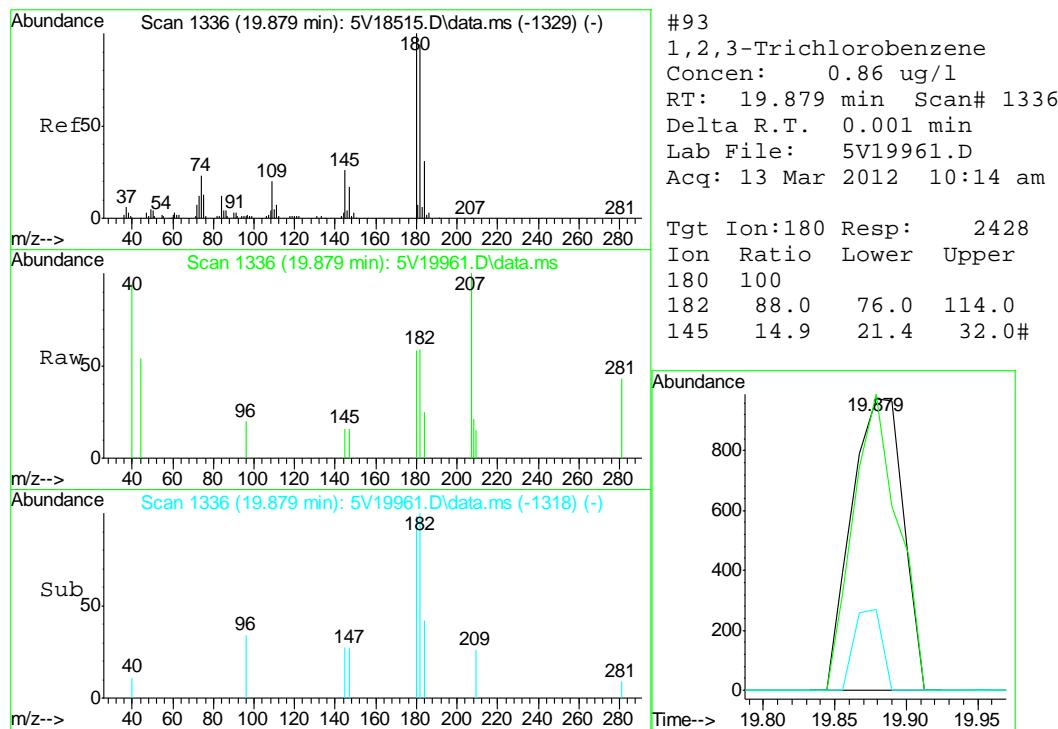
## Quantitation Report (QT Reviewed)

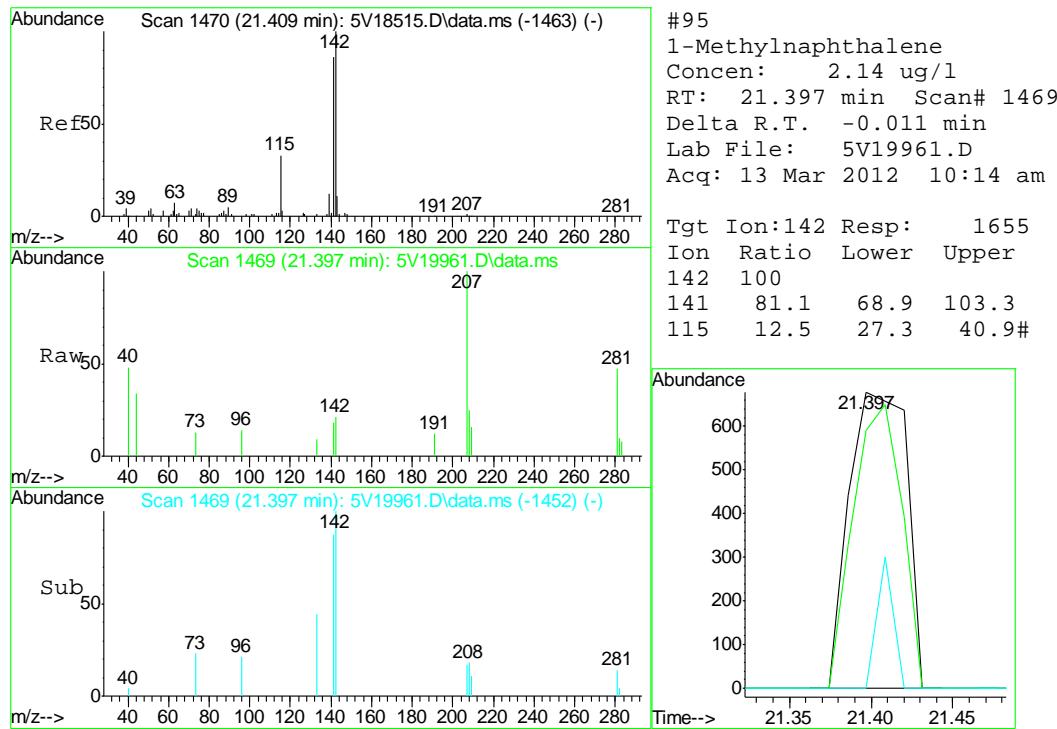
Data Path : C:\msdchem\1\DATA\V5031312.S\  
 Data File : 5V19961.D  
 Acq On : 13 Mar 2012 10:14 am  
 Operator : KOROUSHV  
 Sample : MB  
 Misc : MS3551,V5V1203,5.00,,100,5,1  
 ALS Vial : 3 Sample Multiplier: 1

Quant Time: Mar 14 09:51:19 2012  
 Quant Method : C:\msdchem\1\METHODS\V5AP1186TVH1186.M  
 Quant Title : 8260  
 QLast Update : Fri Mar 02 14:22:16 2012  
 Response via : Initial Calibration











## GC/MS Semi-volatiles

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### QC Data Summaries

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7

Includes the following where applicable:

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

**Method Blank Summary**

**Job Number:** D32609  
**Account:** XTOKWR XTO Energy  
**Project:** FRU 297-32A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
OP5528-MB	3G08524.D	1	03/14/12	DC	03/13/12	OP5528	E3G348

**The QC reported here applies to the following samples:****Method:** SW846 8270C BY SIM

D32609-1

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	
50-32-8	Benzo(a)pyrene	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	
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	78%
321-60-8	2-Fluorobiphenyl	67%
1718-51-0	Terphenyl-d14	79%

## Blank Spike Summary

Page 1 of 1

Job Number: D32609

Account: XTOKWR XTO Energy

Project: FRU 297-32A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
OP5528-BS	3G08525.D	1	03/14/12	DC	03/13/12	OP5528	E3G348

The QC reported here applies to the following samples:

Method: SW846 8270C BY SIM

D32609-1

CAS No.	Compound	Spike ug/kg	BSP ug/kg	BSP %	Limits
83-32-9	Acenaphthene	83.3	66.1	79	34-130
120-12-7	Anthracene	83.3	76.1	91	35-130
56-55-3	Benzo(a)anthracene	83.3	73.1	88	36-130
50-32-8	Benzo(a)pyrene	83.3	69.6	84	36-130
205-99-2	Benzo(b)fluoranthene	83.3	71.1	85	35-130
207-08-9	Benzo(k)fluoranthene	83.3	67.8	81	37-130
218-01-9	Chrysene	83.3	76.8	92	40-130
53-70-3	Dibenzo(a,h)anthracene	83.3	74.9	90	32-130
206-44-0	Fluoranthene	83.3	80.2	96	38-130
86-73-7	Fluorene	83.3	69.8	84	35-130
193-39-5	Indeno(1,2,3-cd)pyrene	83.3	80.9	97	28-130
91-20-3	Naphthalene	83.3	64.5	77	35-130
129-00-0	Pyrene	83.3	72.7	87	37-130

CAS No.	Surrogate Recoveries	BSP	Limits
4165-60-0	Nitrobenzene-d5	76%	10-145%
321-60-8	2-Fluorobiphenyl	66%	10-130%
1718-51-0	Terphenyl-d14	78%	22-130%

# Matrix Spike/Matrix Spike Duplicate Summary

Page 1 of 1

Job Number: D32609

Account: XTOKWR XTO Energy

Project: FRU 297-32A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
OP5528-MS	3G08527.D	1	03/14/12	DC	03/13/12	OP5528	E3G348
OP5528-MSD	3G08528.D	1	03/14/12	DC	03/13/12	OP5528	E3G348
D32609-1	3G08526.D	1	03/14/12	DC	03/13/12	OP5528	E3G348

The QC reported here applies to the following samples:

Method: SW846 8270C BY SIM

D32609-1

CAS No.	Compound	D32609-1		Spike	MS	MS	MSD	MSD	RPD	Limits Rec/RPD
		ug/kg	Q	ug/kg	ug/kg	%	ug/kg	%		
83-32-9	Acenaphthene	ND	88	43.1	49	61.7	70	35* a	10-155/30	
120-12-7	Anthracene	ND	88	53.4	61	70.2	80	27	10-155/30	
56-55-3	Benzo(a)anthracene	ND	88	52.8	60	70.4	80	29	10-175/30	
50-32-8	Benzo(a)pyrene	ND	88	45.2	51	60.2	68	28	10-164/30	
205-99-2	Benzo(b)fluoranthene	ND	88	48.8	55	62.7	71	25	10-165/30	
207-08-9	Benzo(k)fluoranthene	ND	88	40.6	46	50.3	57	21	10-178/30	
218-01-9	Chrysene	ND	88	51.8	59	71.4	81	32* a	10-147/30	
53-70-3	Dibenzo(a,h)anthracene	ND	88	74.1	84	108	123	37* a	10-144/30	
206-44-0	Fluoranthene	ND	88	50.5	57	65.2	74	25	10-207/30	
86-73-7	Fluorene	ND	88	56.7	64	79.0	90	33* a	10-163/30	
193-39-5	Indeno(1,2,3-cd)pyrene	ND	88	83.3	95	119	135	35* a	10-180/30	
91-20-3	Naphthalene	32.7	88	84.7	59	125	105	38* a	10-198/30	
129-00-0	Pyrene	ND	88	60.4	69	81.5	93	30	10-189/30	

CAS No.	Surrogate Recoveries	MS	MSD	D32609-1	Limits
4165-60-0	Nitrobenzene-d5	49%	70%	53%	10-145%
321-60-8	2-Fluorobiphenyl	40%	59%	48%	10-130%
1718-51-0	Terphenyl-d14	57%	77%	76%	22-130%

(a) Variability of recovery may be due to sample matrix/homogeneity.

7.3.1  
7



## GC/MS Semi-volatiles

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Raw Data

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## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\031412\  
 Data File : 3g08526.D  
 Acq On : 14 Mar 2012 10:49 am  
 Operator : DONC  
 Sample : D32609-1  
 Misc : OP5528,E3G348,30.11,,,1,1  
 ALS Vial : 6 Sample Multiplier: 1

Quant Time: Mar 15 08:02:09 2012  
 Quant Method : C:\msdchem\1\METHODS\SIMPE3G344.M  
 Quant Title : PAHSIM BASE  
 QLast Update : Mon Mar 12 09:19:25 2012  
 Response via : Initial Calibration

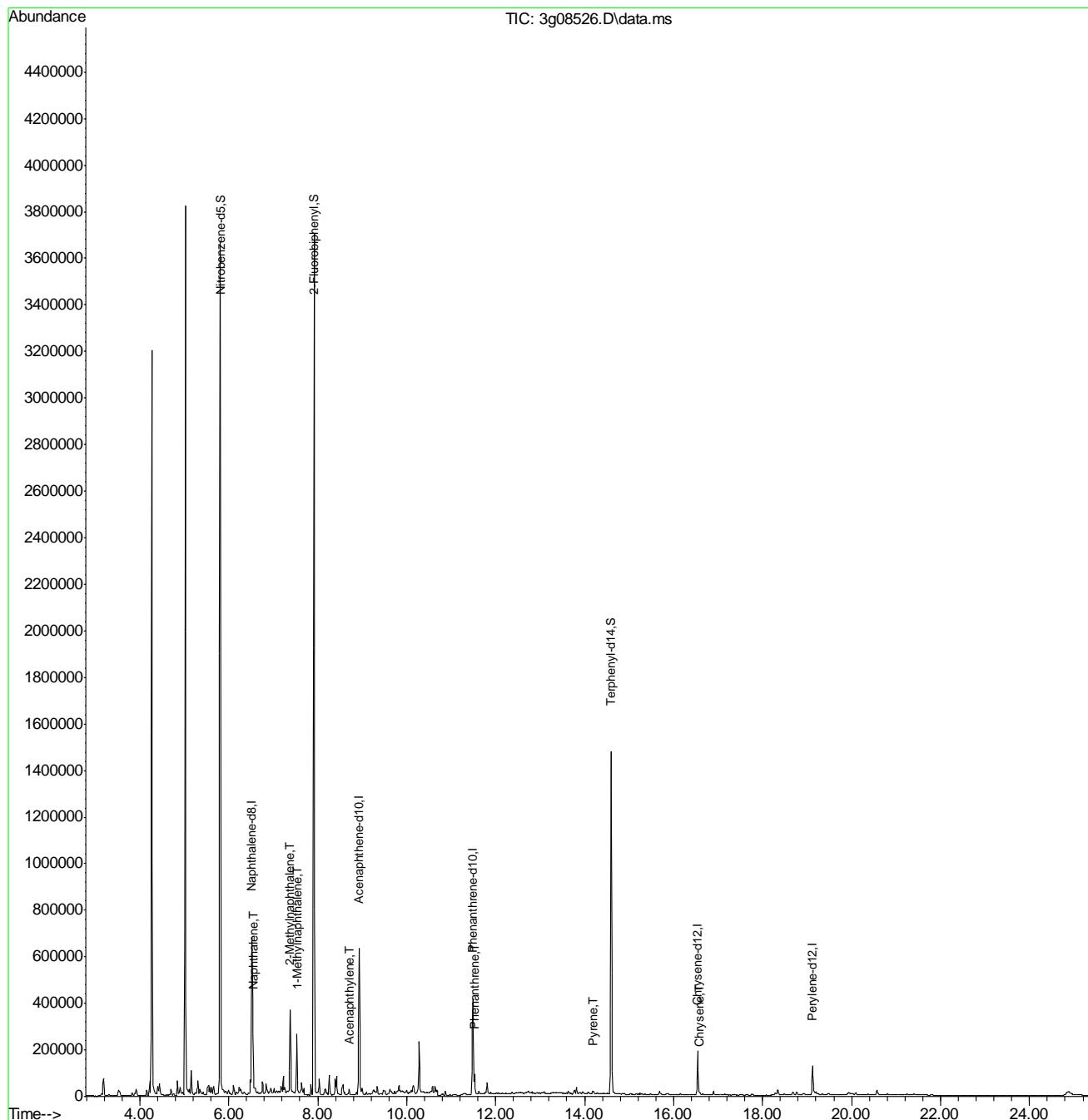
Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
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Internal Standards						
1) Naphthalene-d8	6.520	136	626276	4.00	ug/mL	-0.01
6) Acenaphthene-d10	8.933	164	326715	4.00	ug/mL	0.00
14) Phenanthrene-d10	11.485	188	411909	4.00	ug/mL	0.00
18) Chrysene-d12	16.540	240	210352	4.00	ug/mL	-0.01
23) Perylene-d12	19.122	264	172963	4.00	ug/mL	-0.01
<hr/>						
System Monitoring Compounds						
2) Nitrobenzene-d5	5.809	82	2244963	26.30	ug/mL	0.00
Spiked Amount 50.000	Range 25 - 135		Recovery	=	52.60%	
7) 2-Fluorobiphenyl	7.917	172	3155296	24.07	ug/mL	-0.01
Spiked Amount 50.000	Range 25 - 135		Recovery	=	48.14%	
20) Terphenyl-d14	14.595	244	1726107	37.96	ug/mL	-0.02
Spiked Amount 50.000	Range 25 - 135		Recovery	=	75.92%	
<hr/>						
Target Compounds						
3) N-Nitrosodimethylamine	0.000		0	N.D.	d	
4) N-Nitrosodi-propylamine	0.000		0	N.D.	d	
5) Naphthalene	6.545	128	195653	0.93	ug/mL	98
8) 2-Methylnaphthalene	7.380	142	196965	1.66	ug/mL	99
9) 1-Methylnaphthalene	7.530	142	102075	0.90	ug/mL	99
10) Acenaphthylene	8.708	152	3313	0.02	ug/mL#	1
11) Acenaphthene	0.000		0	N.D.	d	
12) Fluorene	0.000		0	N.D.	d	
13) Diphenylamine	0.000		0	N.D.	d	
15) Phenanthrene	11.524	178	95917	0.67	ug/mL	96
16) Anthracene	0.000		0	N.D.	d	
17) Fluoranthene	0.000		0	N.D.	d	
19) Pyrene	14.184	202	6974	0.07	ug/mL#	1
21) Benzo(a)anthracene	0.000		0	N.D.	d	
22) Chrysene	16.587	228	8504	0.12	ug/mL	92
24) Benzo(b)fluoranthene	0.000		0	N.D.	d	
25) Benzo(k)fluoranthene	0.000		0	N.D.	d	
26) Benzo(a)pyrene	0.000		0	N.D.	d	
27) Indeno(1,2,3-cd)pyrene	0.000		0	N.D.	d	
28) Dibenz(a,h)anthracene	0.000		0	N.D.	d	
29) Benzo(g,h,i)perylene	0.000		0	N.D.	d	

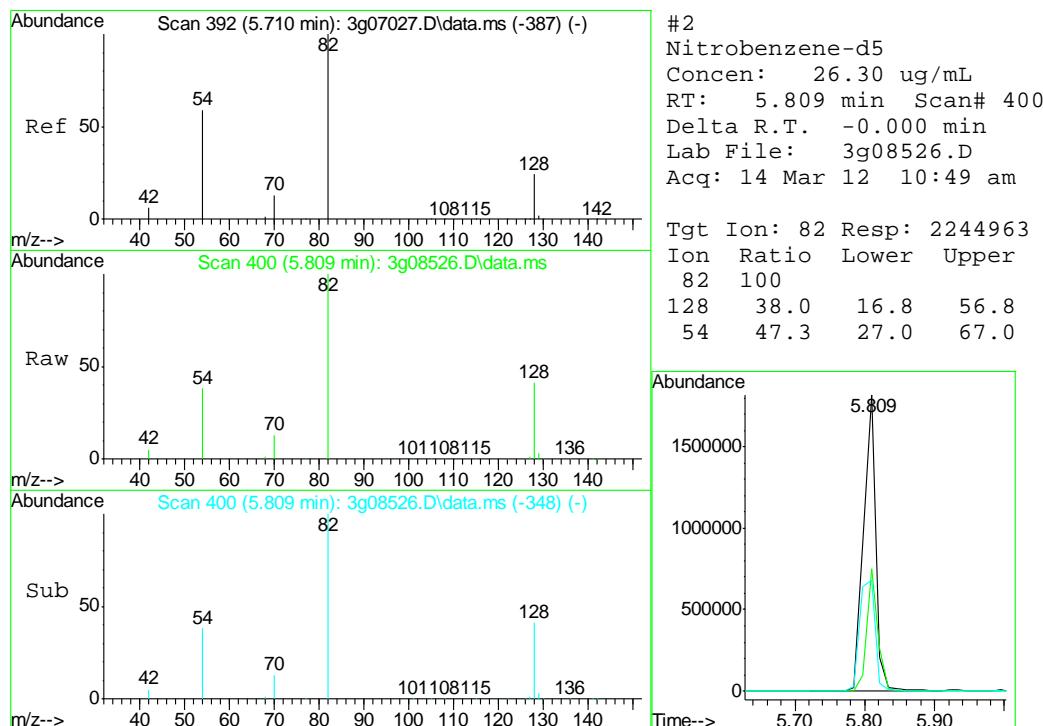
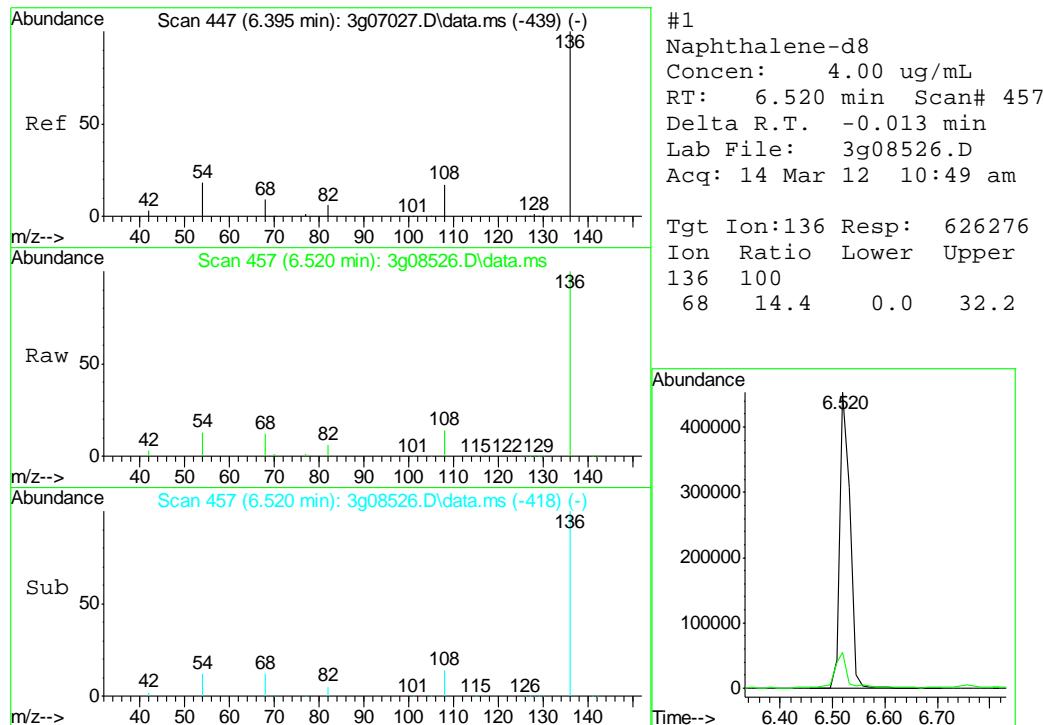
(#) = qualifier out of range (m) = manual integration (+) = signals summed

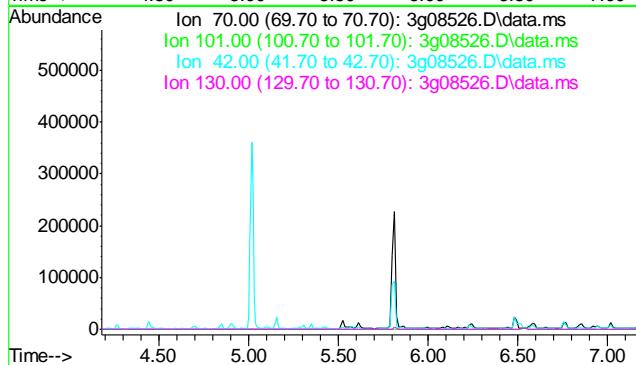
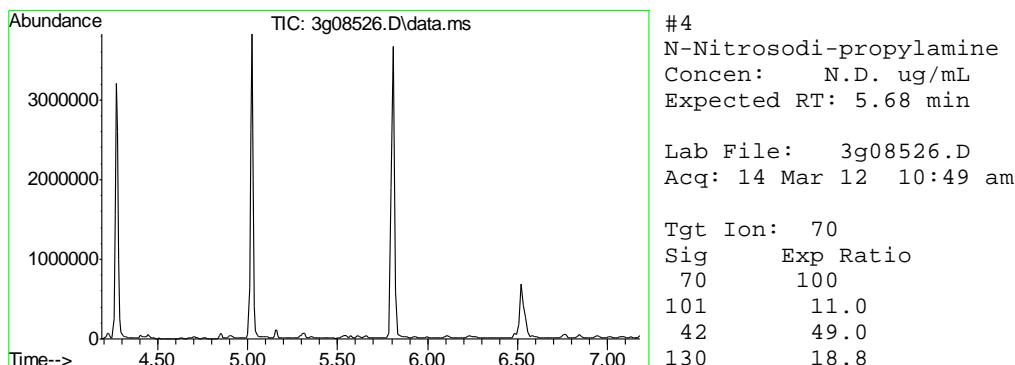
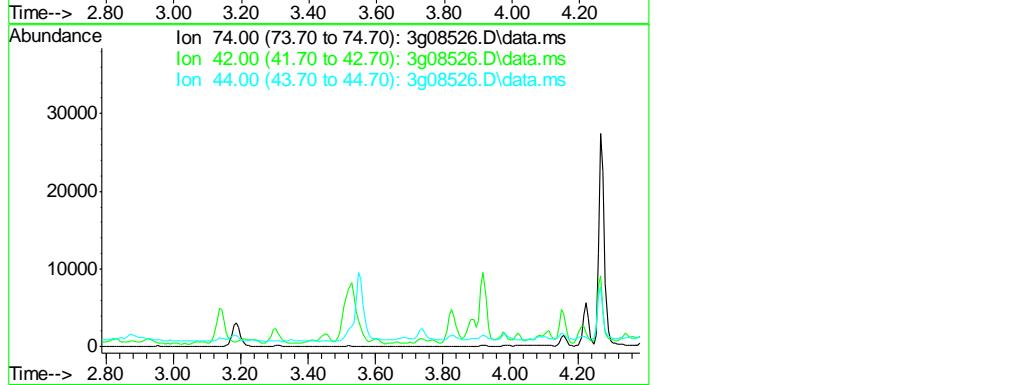
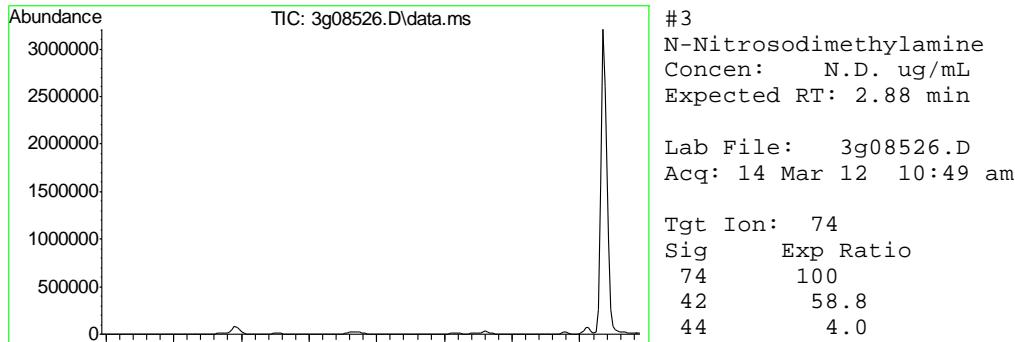
## Quantitation Report (QT Reviewed)

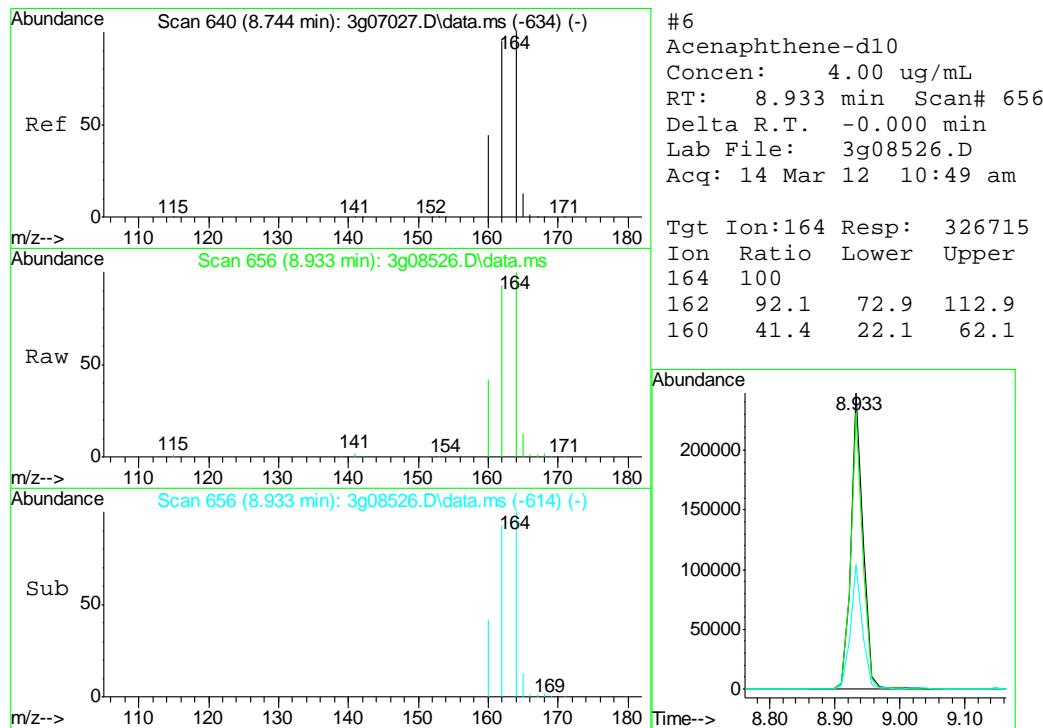
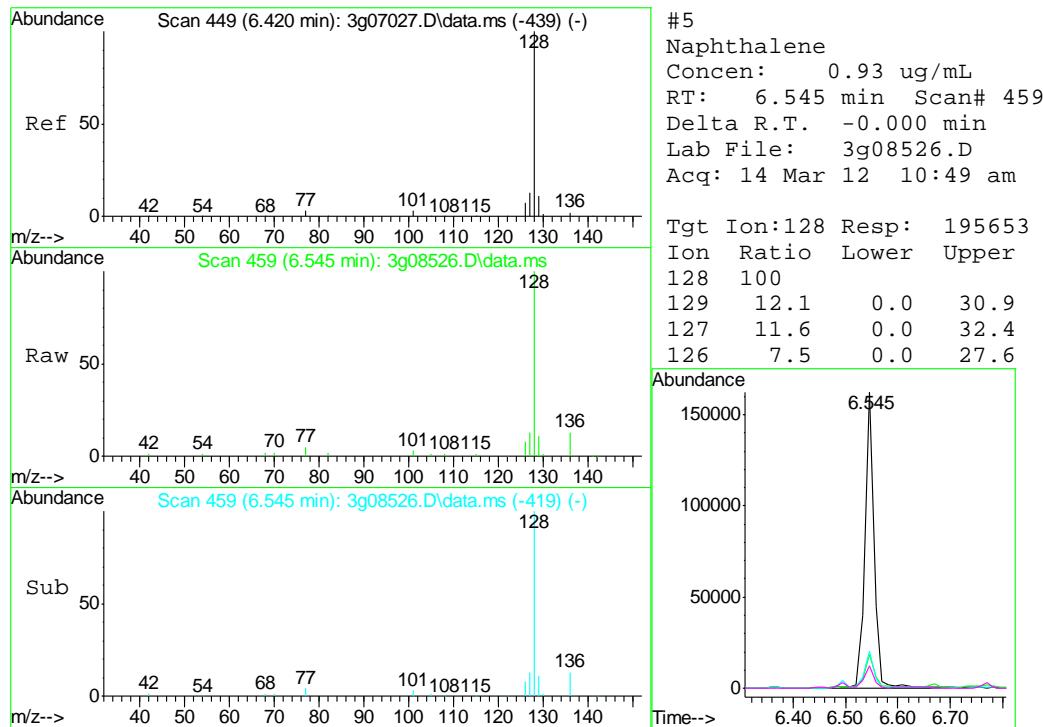
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 Data File : 3g08526.D  
 Acq On : 14 Mar 2012 10:49 am  
 Operator : DONC  
 Sample : D32609-1  
 Misc : OP5528,E3G348,30.11,,,1,1  
 ALS Vial : 6 Sample Multiplier: 1

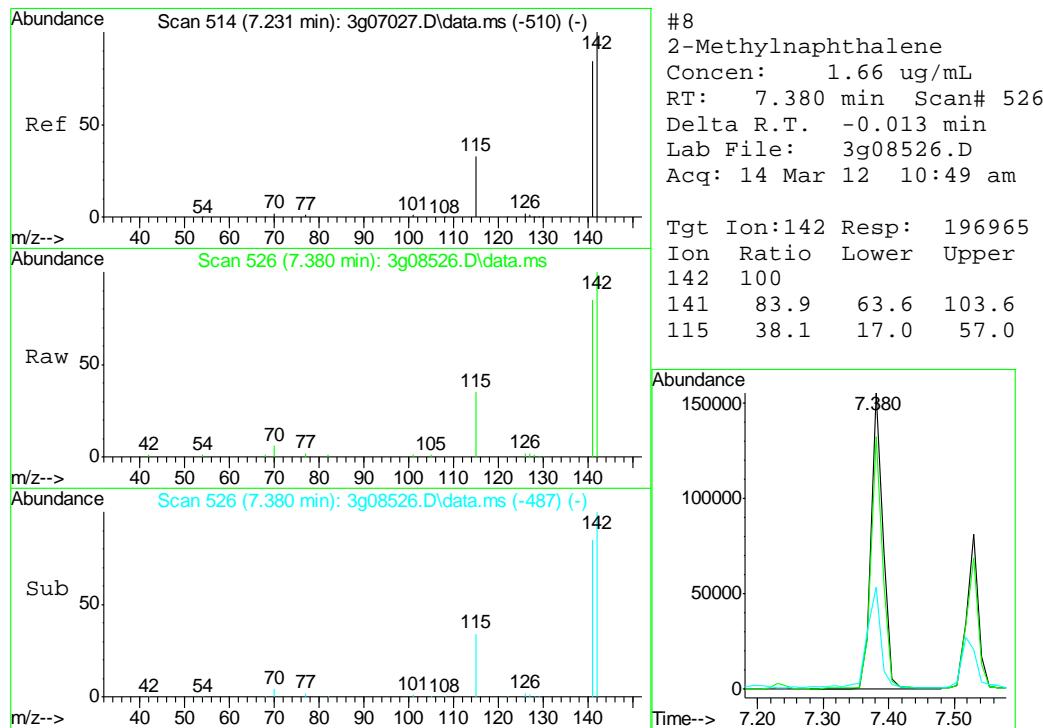
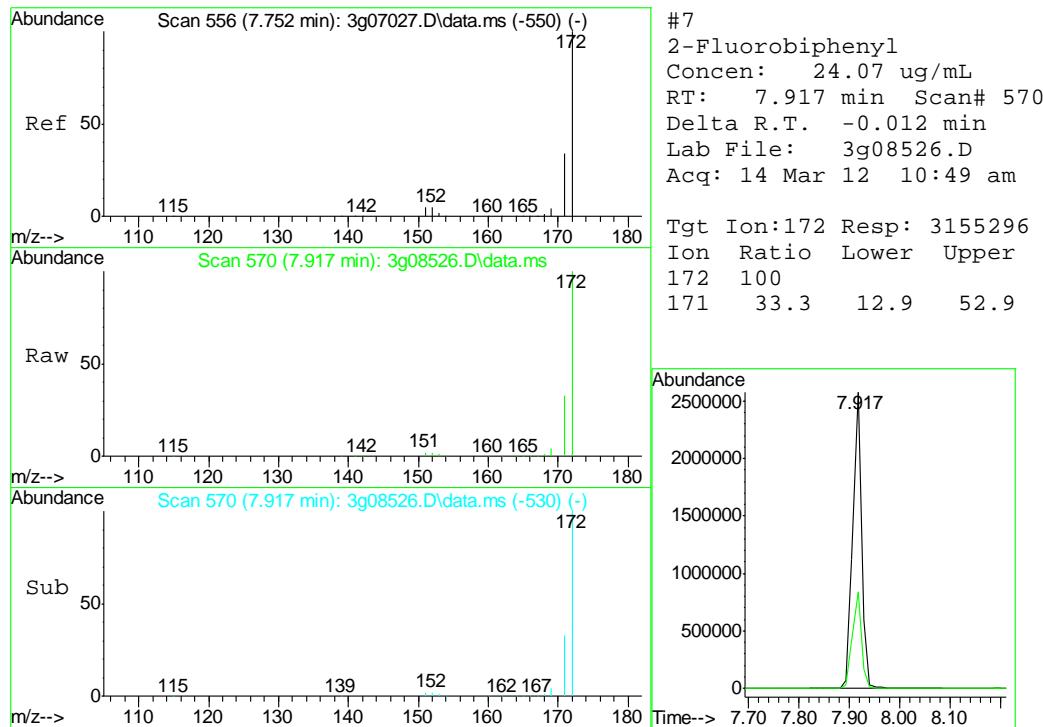
Quant Time: Mar 15 08:02:09 2012  
 Quant Method : C:\msdchem\1\METHODS\SIMPE3G344.M  
 Quant Title : PAHSIM BASE  
 QLast Update : Mon Mar 12 09:19:25 2012  
 Response via : Initial Calibration

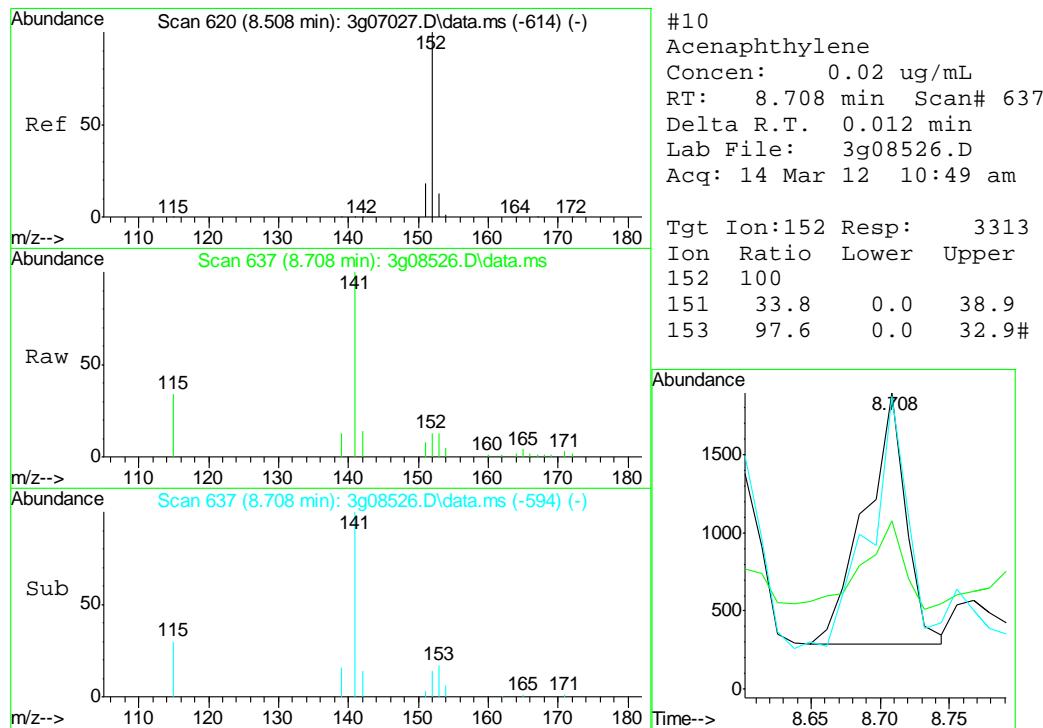
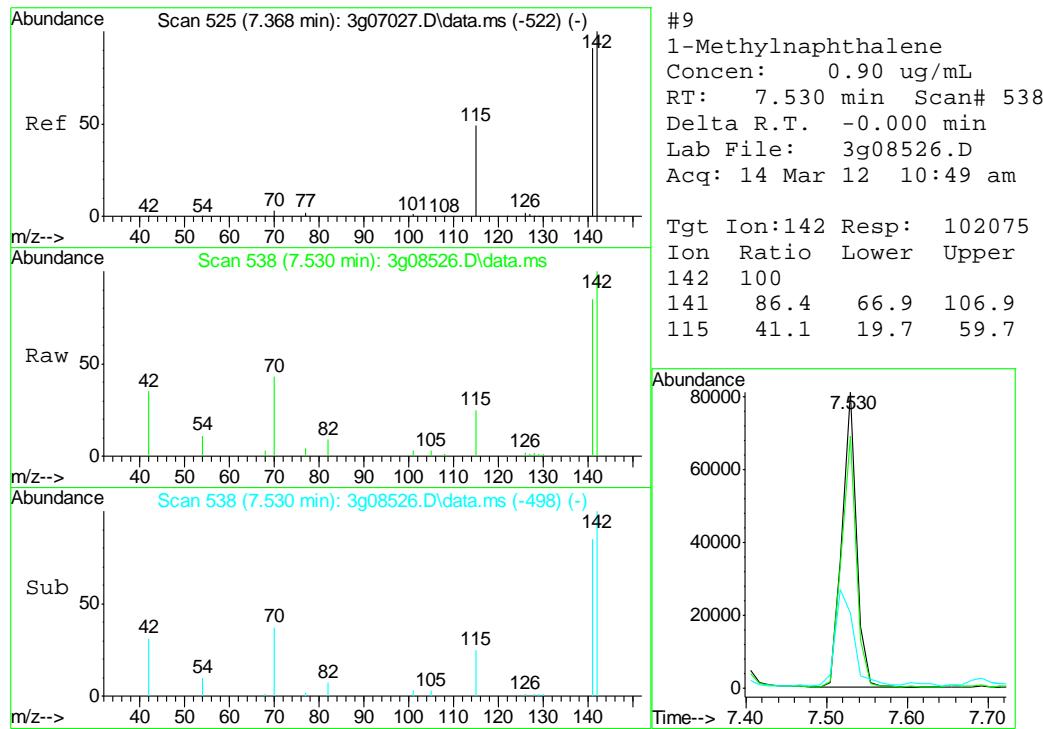


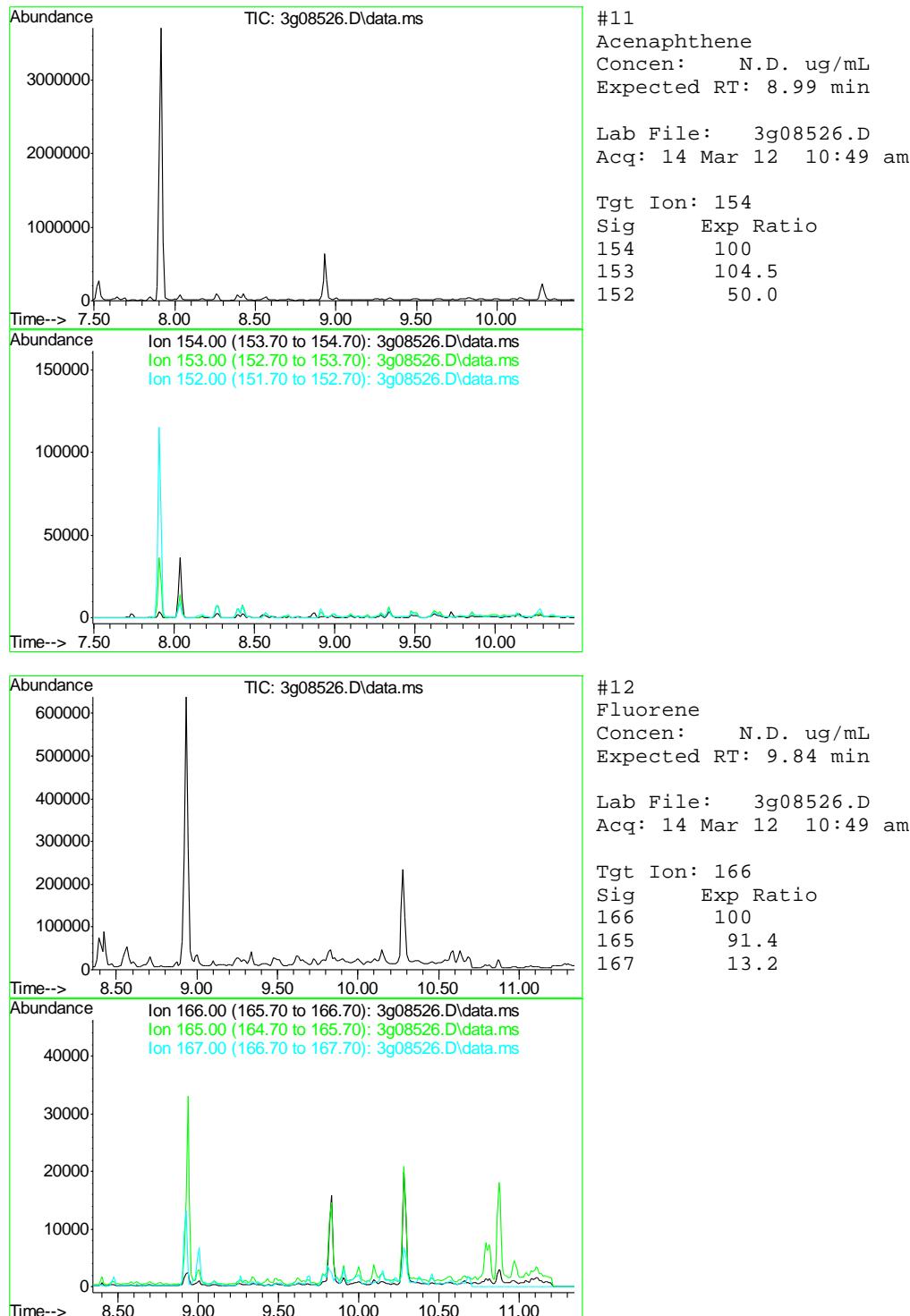


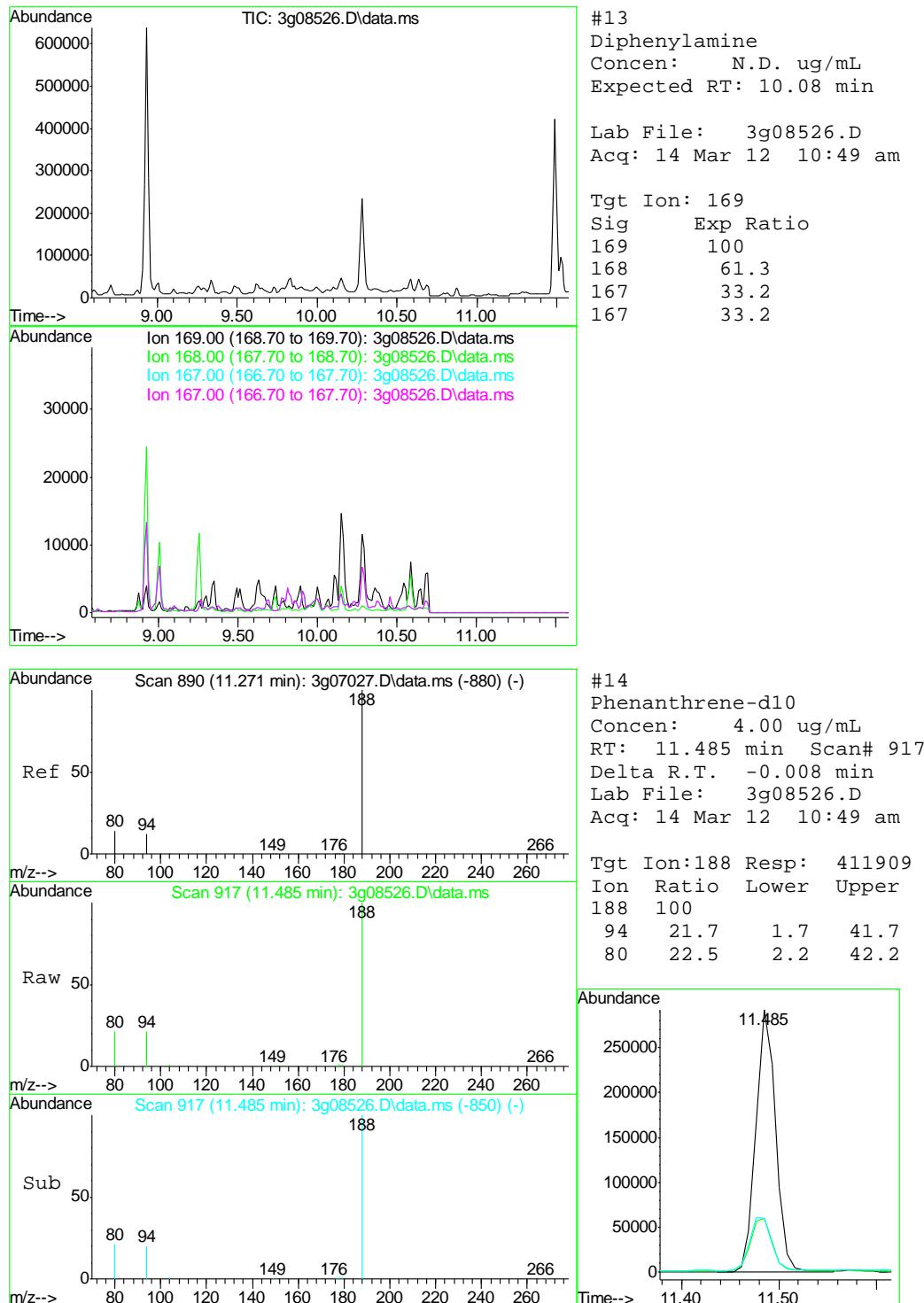


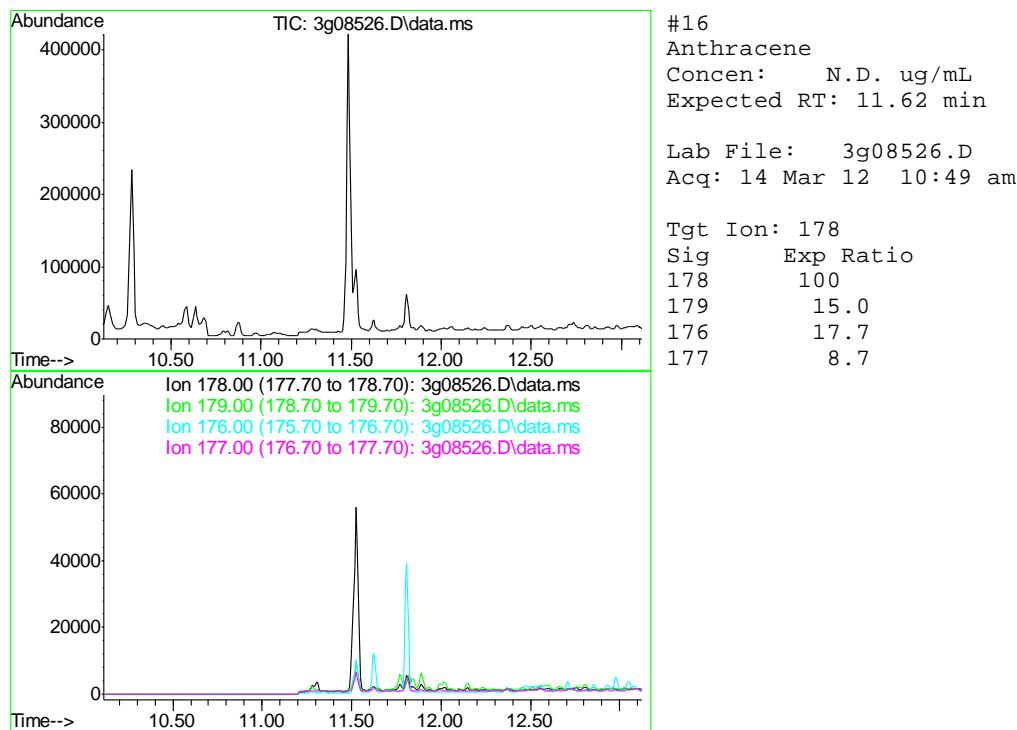
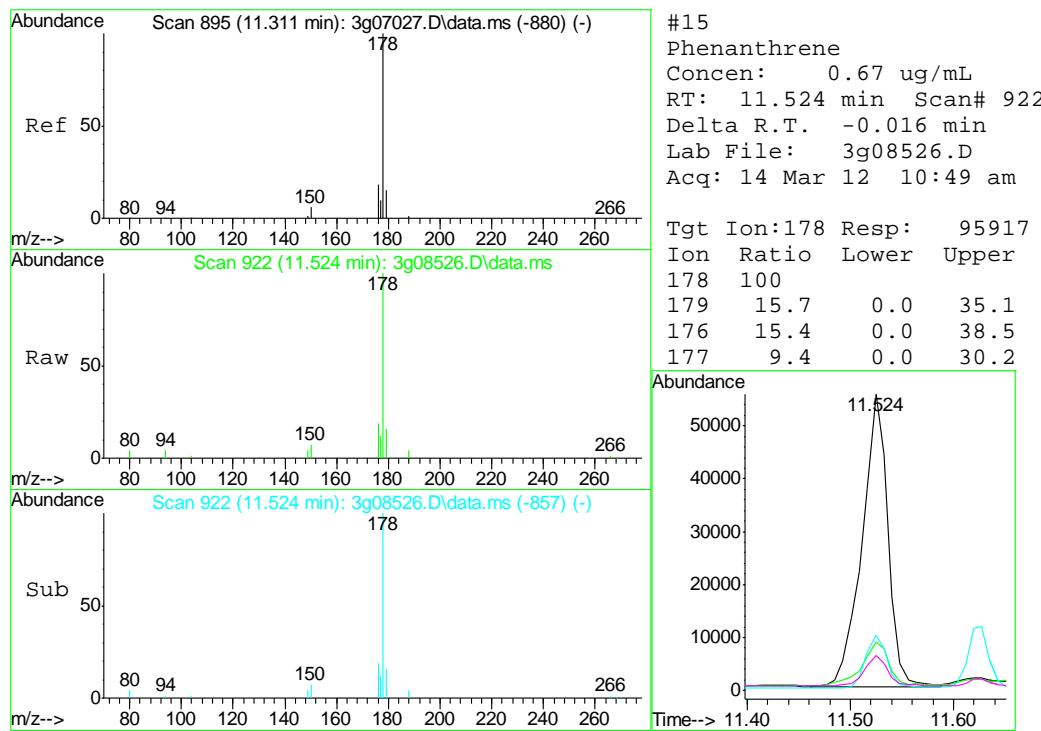


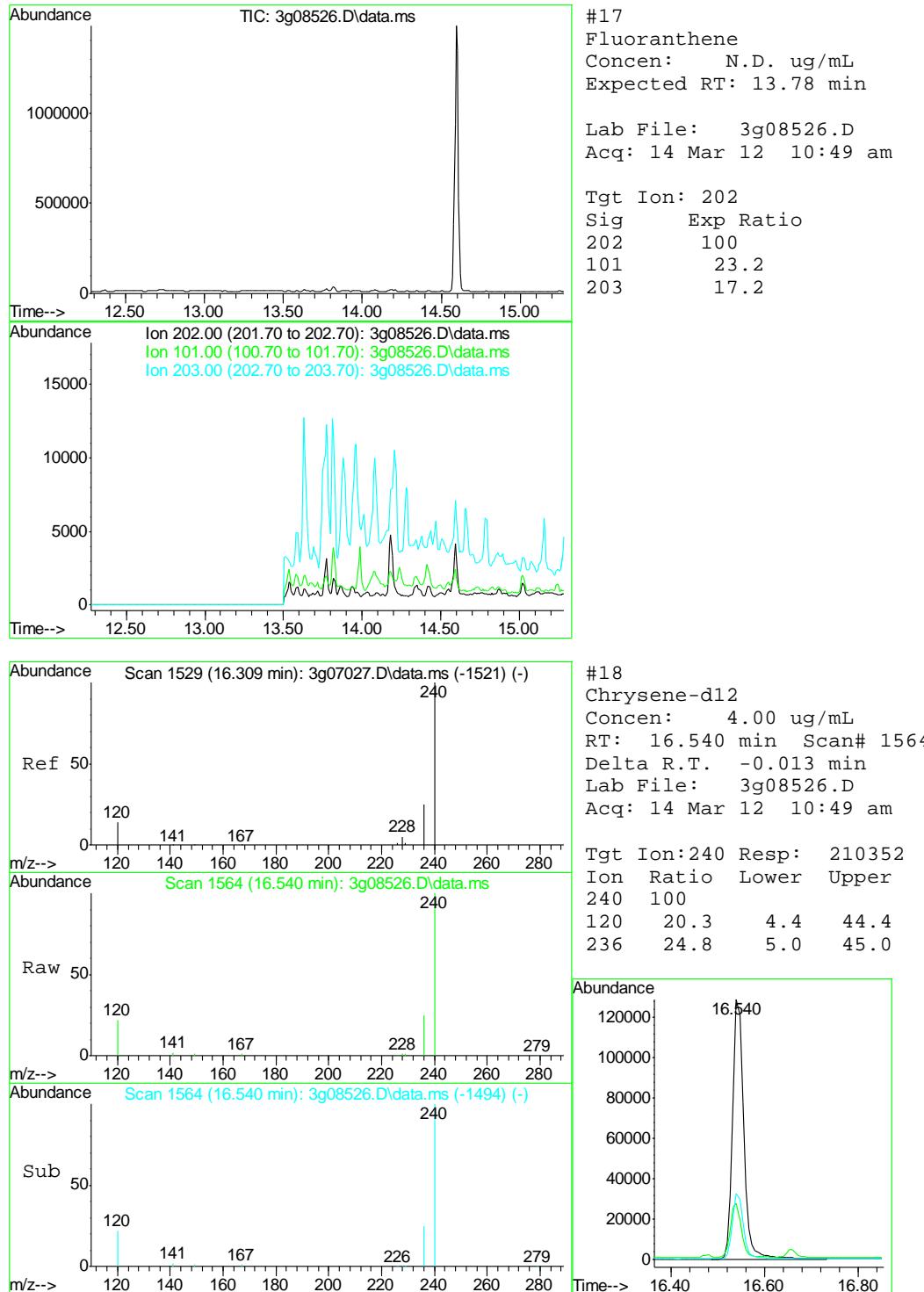


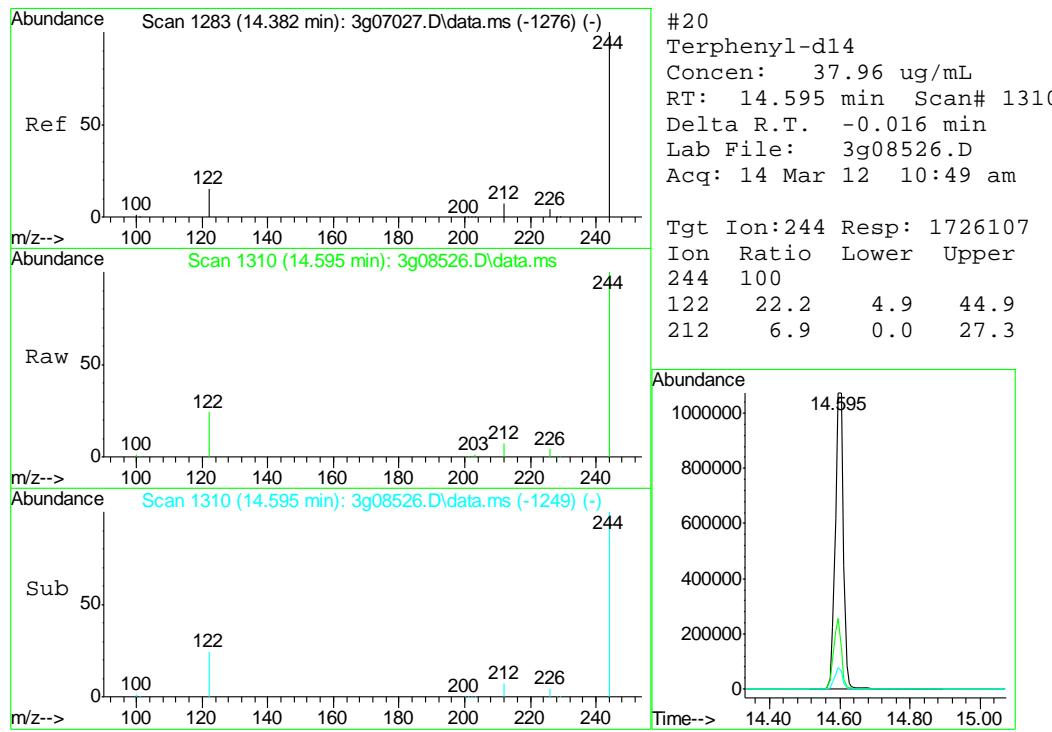
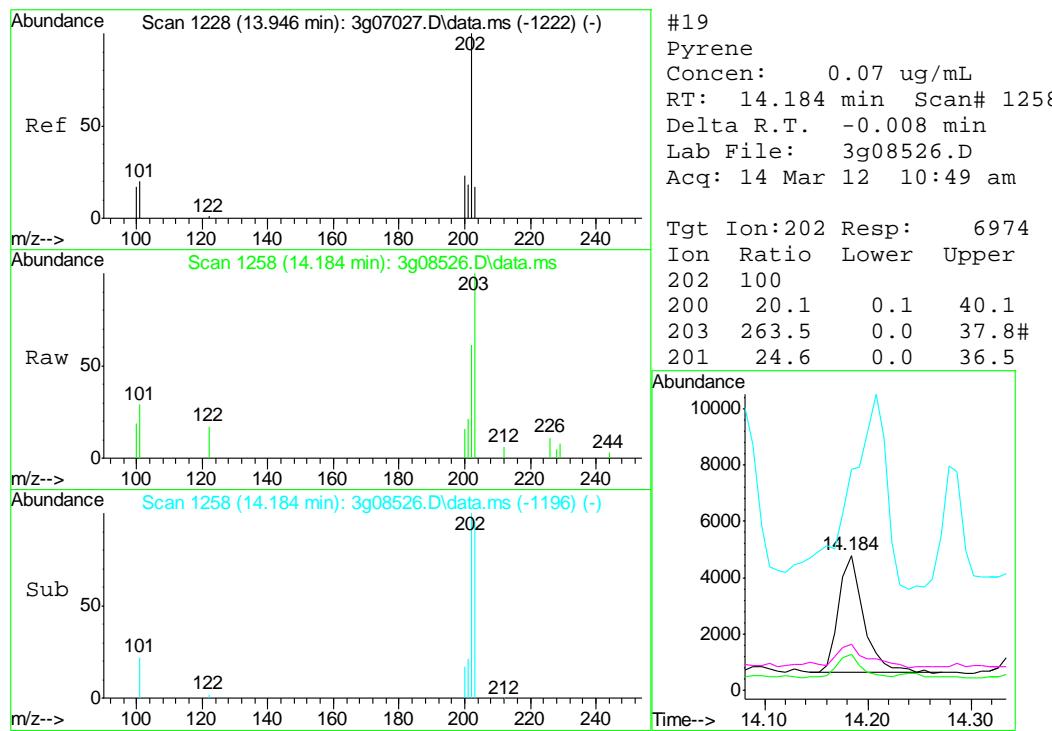


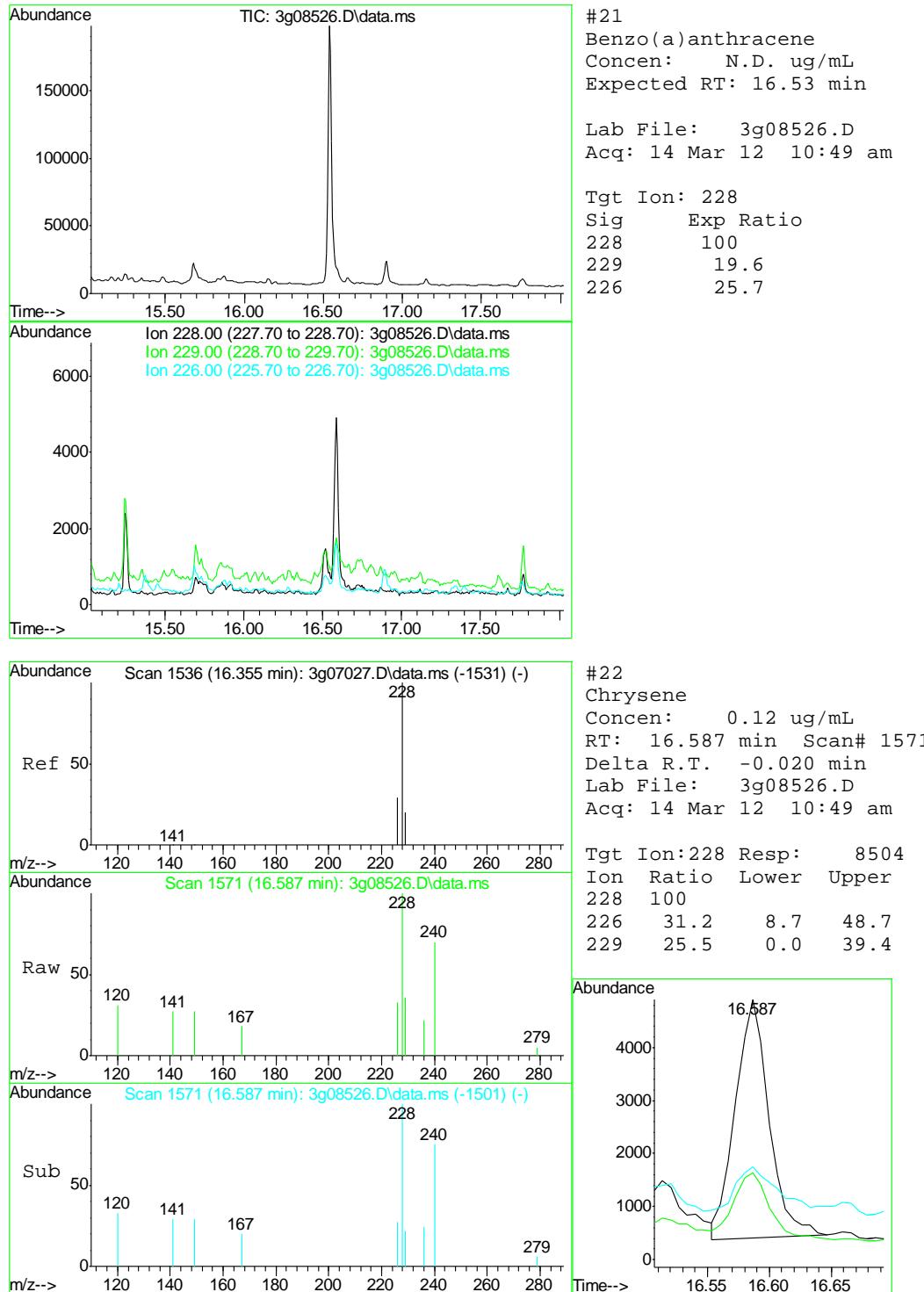


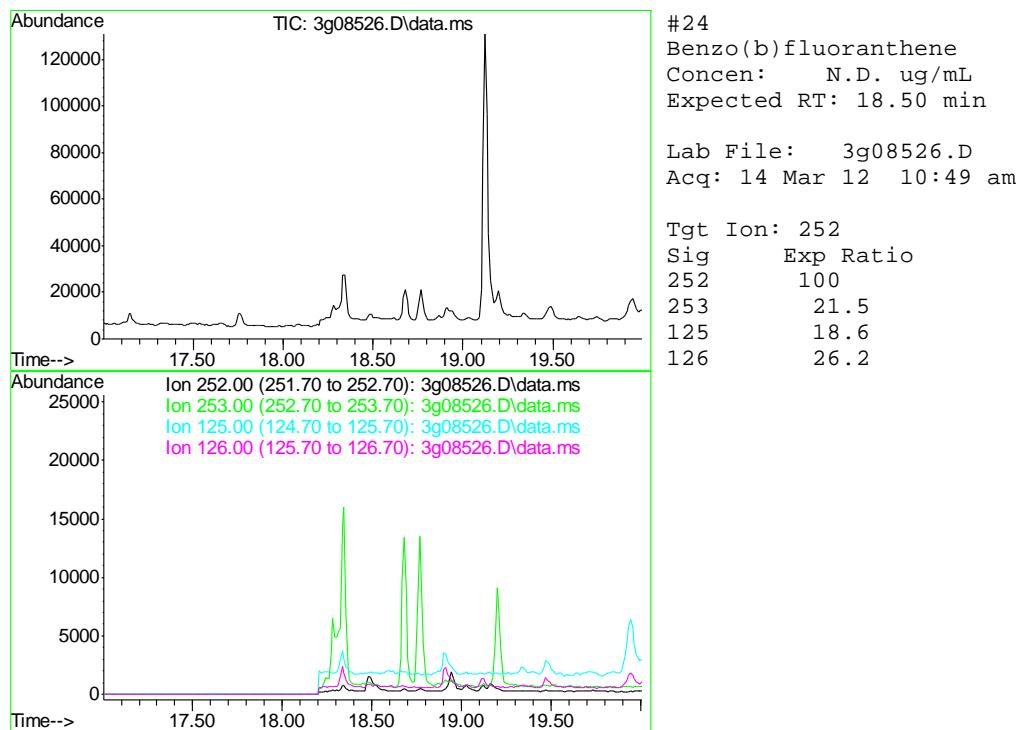
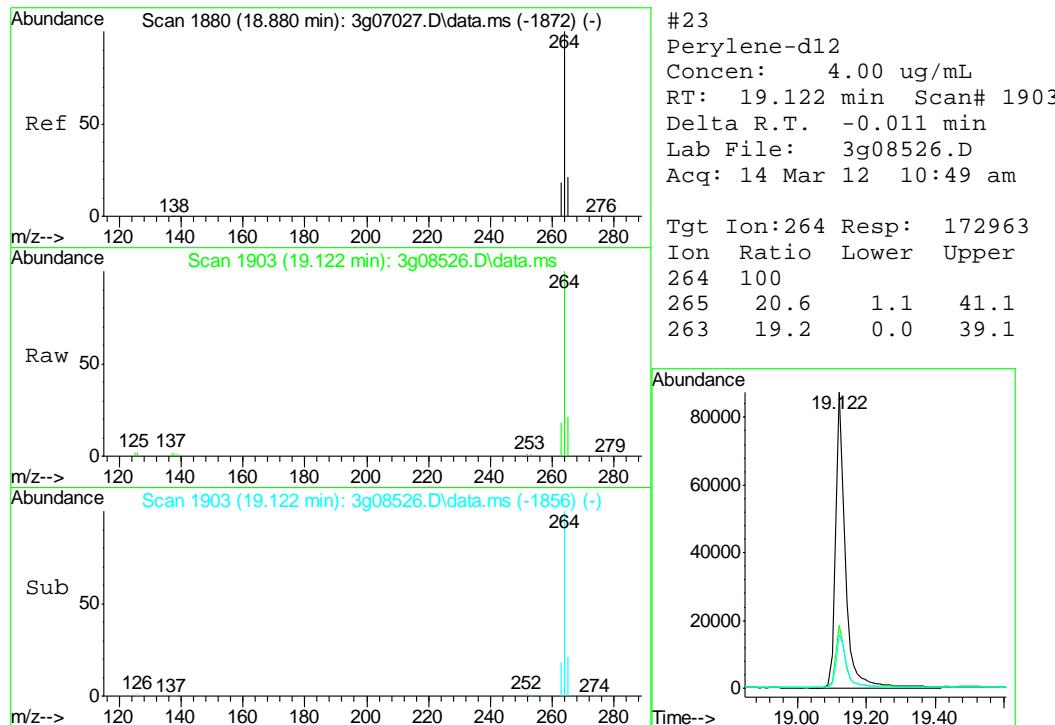


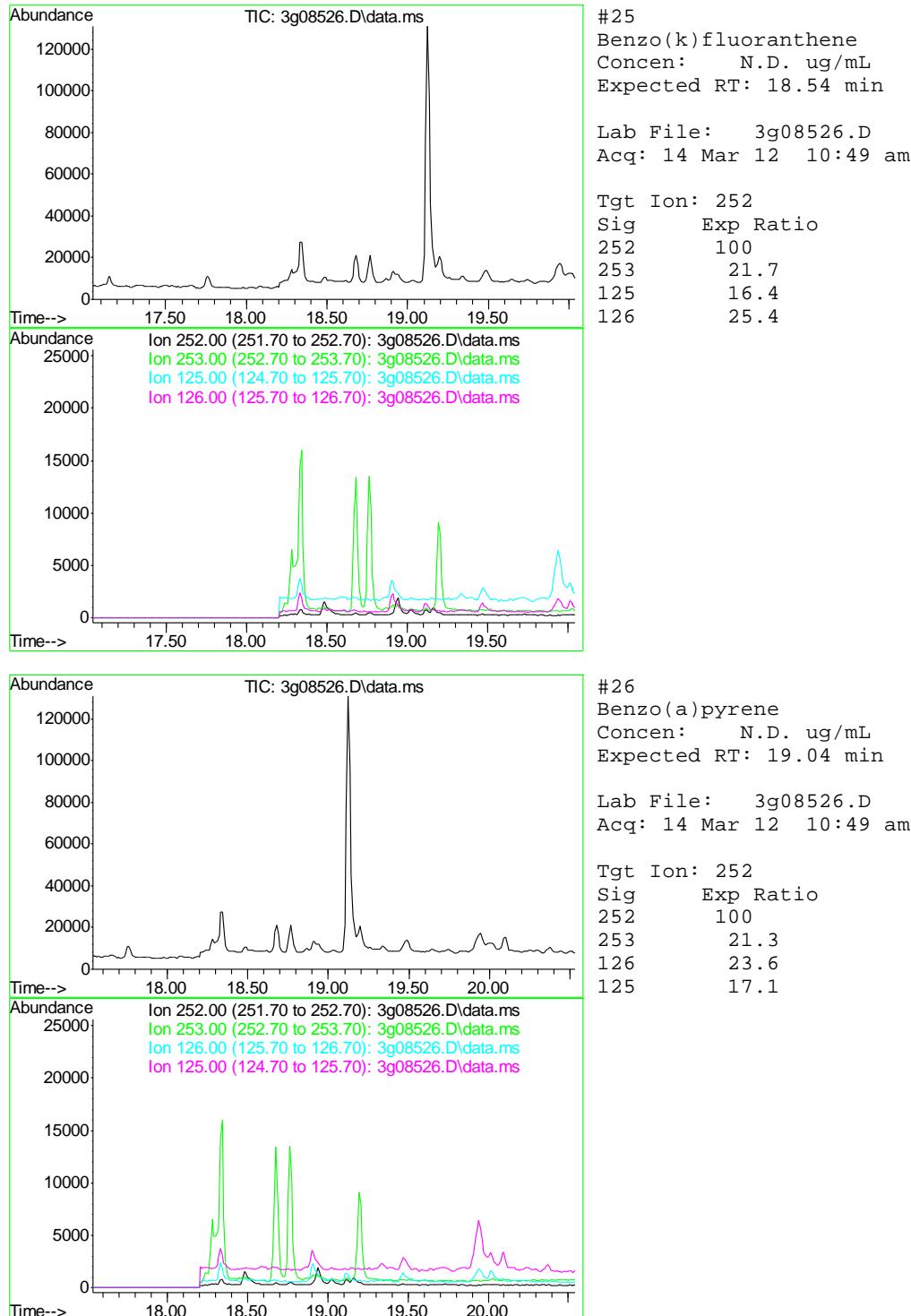


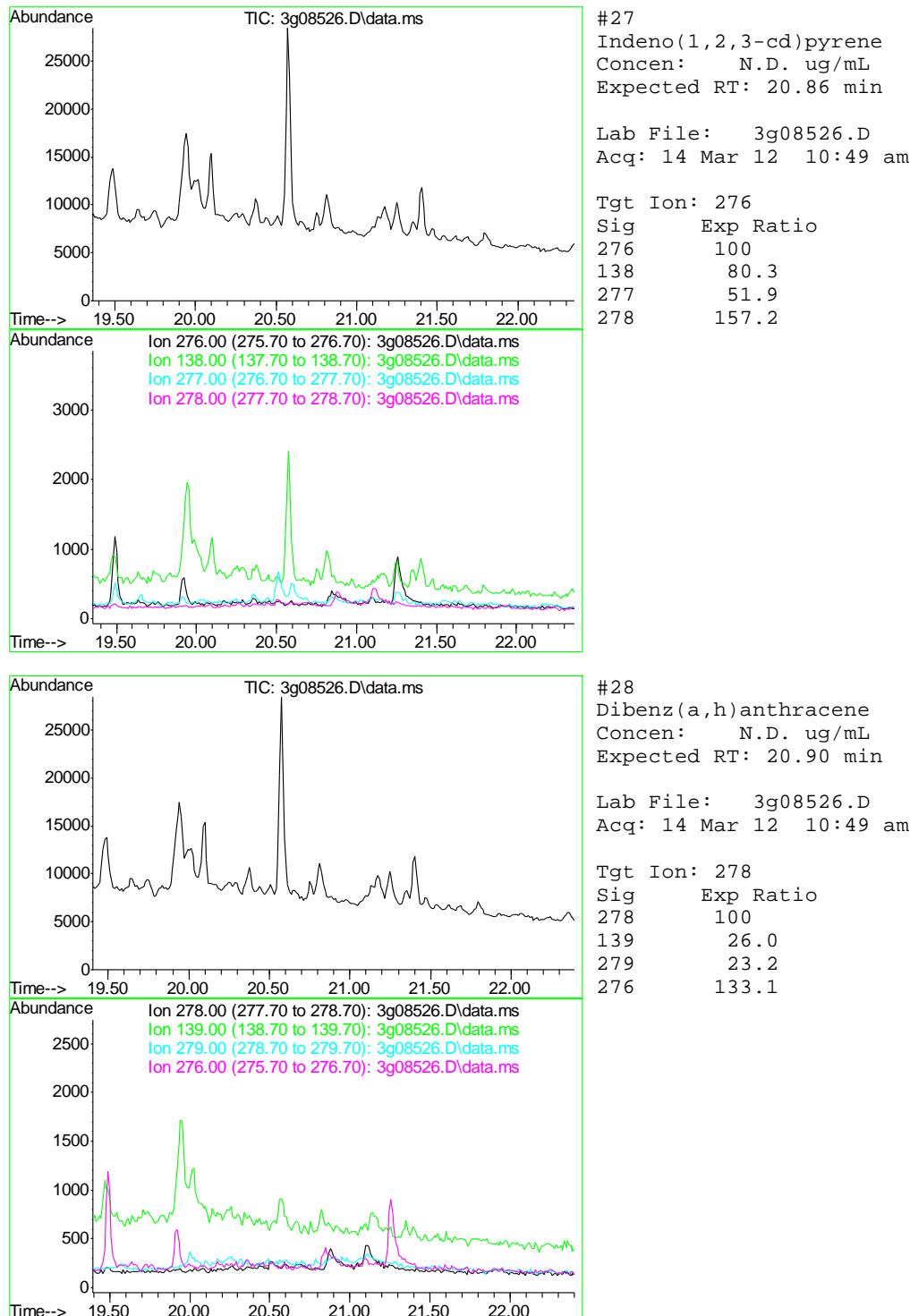


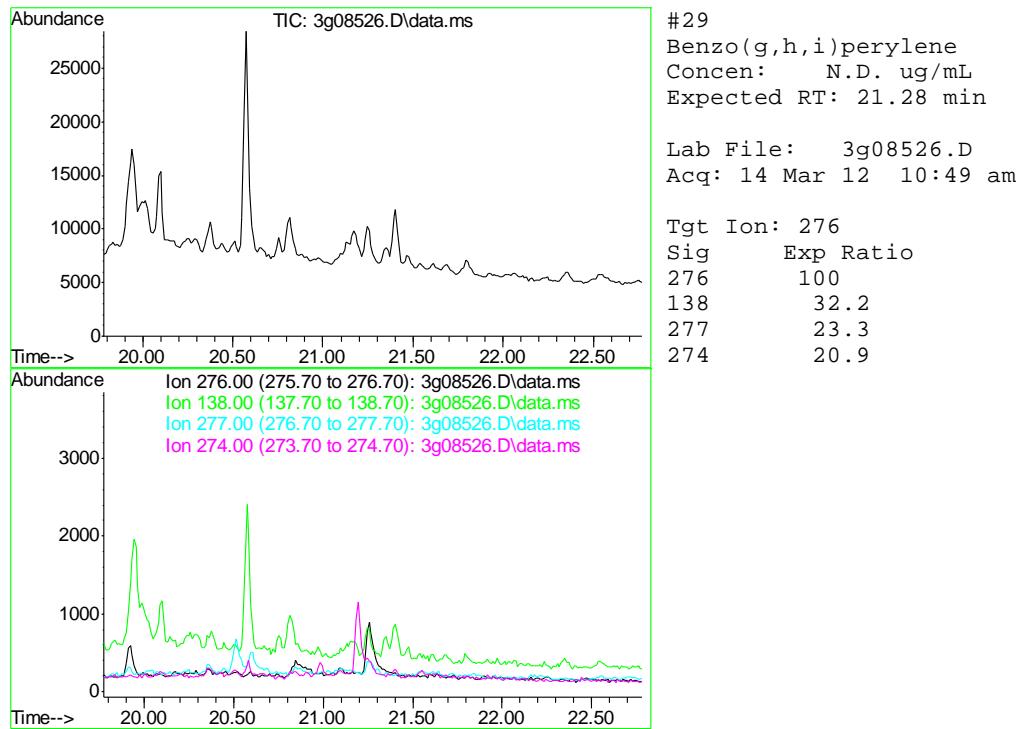












## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\031412\  
 Data File : 3g08524.D  
 Acq On : 14 Mar 2012 9:39 am  
 Operator : DONC  
 Sample : OP5528-MB  
 Misc : OP5528,E3G348,30.00,,,1,1  
 ALS Vial : 4 Sample Multiplier: 1

Quant Time: Mar 15 07:59:40 2012  
 Quant Method : C:\msdchem\1\METHODS\SIMPE3G344.M  
 Quant Title : PAHSIM BASE  
 QLast Update : Mon Mar 12 09:19:25 2012  
 Response via : Initial Calibration

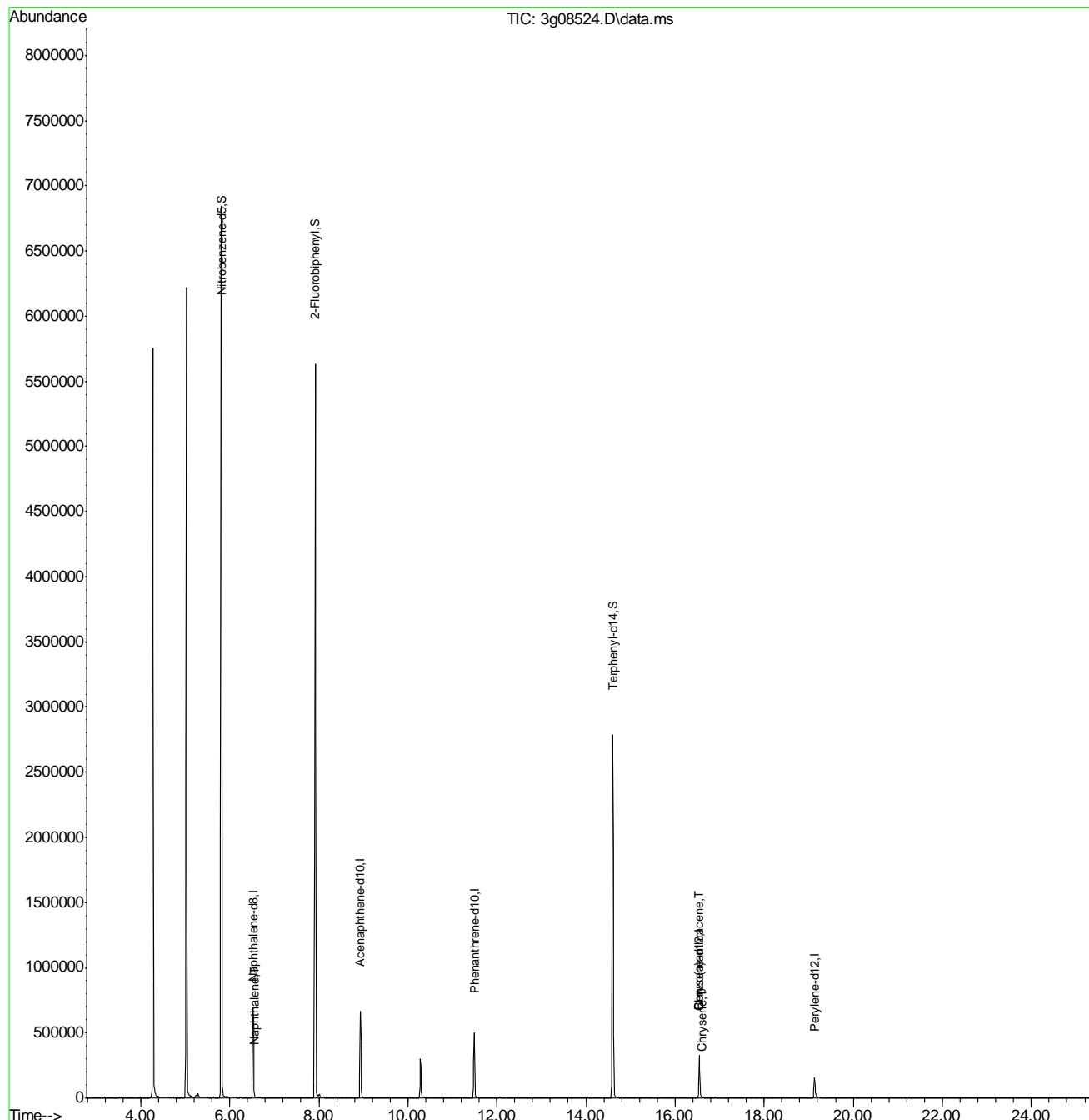
Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
<hr/>						
Internal Standards						
1) Naphthalene-d8	6.532	136	745037	4.00	ug/mL	0.00
6) Acenaphthene-d10	8.933	164	381870	4.00	ug/mL	0.00
14) Phenanthrene-d10	11.493	188	534122	4.00	ug/mL	0.00
18) Chrysene-d12	16.547	240	397538	4.00	ug/mL	0.00
23) Perylene-d12	19.132	264	263103	4.00	ug/mL	0.00
<hr/>						
System Monitoring Compounds						
2) Nitrobenzene-d5	5.809	82	3959513	39.00	ug/mL	0.00
Spiked Amount 50.000	Range 25 - 135		Recovery	=	78.00%	
7) 2-Fluorobiphenyl	7.917	172	5119221	33.41	ug/mL	-0.01
Spiked Amount 50.000	Range 25 - 135		Recovery	=	66.82%	
20) Terphenyl-d14	14.603	244	3392724	39.48	ug/mL	0.00
Spiked Amount 50.000	Range 25 - 135		Recovery	=	78.96%	
<hr/>						
Target Compounds						
3) N-Nitrosodimethylamine	0.000		0	N.D.	d	
4) N-Nitrosodi-propylamine	0.000		0	N.D.	d	
5) Naphthalene	6.545	128	2724	0.01	ug/mL	93
8) 2-Methylnaphthalene	0.000		0	N.D.	d	
9) 1-Methylnaphthalene	0.000		0	N.D.	d	
10) Acenaphthylene	0.000		0	N.D.	d	
11) Acenaphthene	0.000		0	N.D.	d	
12) Fluorene	0.000		0	N.D.		
13) Diphenylamine	0.000		0	N.D.	d	
15) Phenanthrene	0.000		0	N.D.	d	
16) Anthracene	0.000		0	N.D.	d	
17) Fluoranthene	0.000		0	N.D.	d	
19) Pyrene	0.000		0	N.D.	d	
21) Benzo(a)anthracene	16.540	228	1664	0.01	ug/mL	70
22) Chrysene	16.593	228	1201	0.01	ug/mL	82
24) Benzo(b)fluoranthene	0.000		0	N.D.	d	
25) Benzo(k)fluoranthene	0.000		0	N.D.	d	
26) Benzo(a)pyrene	0.000		0	N.D.	d	
27) Indeno(1,2,3-cd)pyrene	0.000		0	N.D.	d	
28) Dibenz(a,h)anthracene	0.000		0	N.D.	d	
29) Benzo(g,h,i)perylene	0.000		0	N.D.	d	

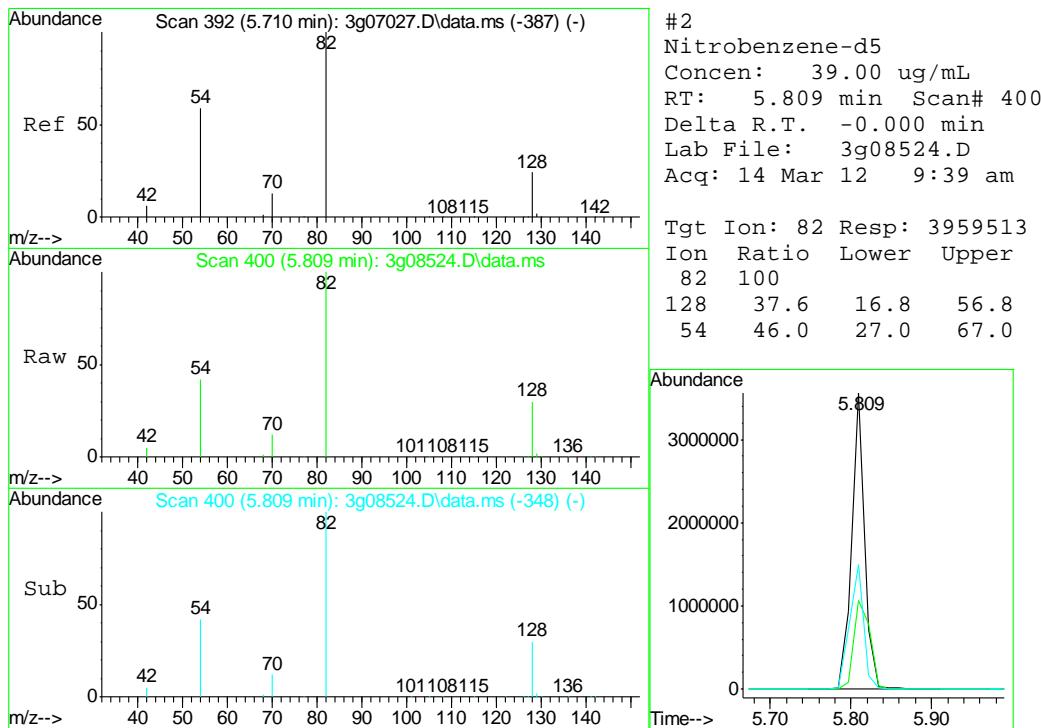
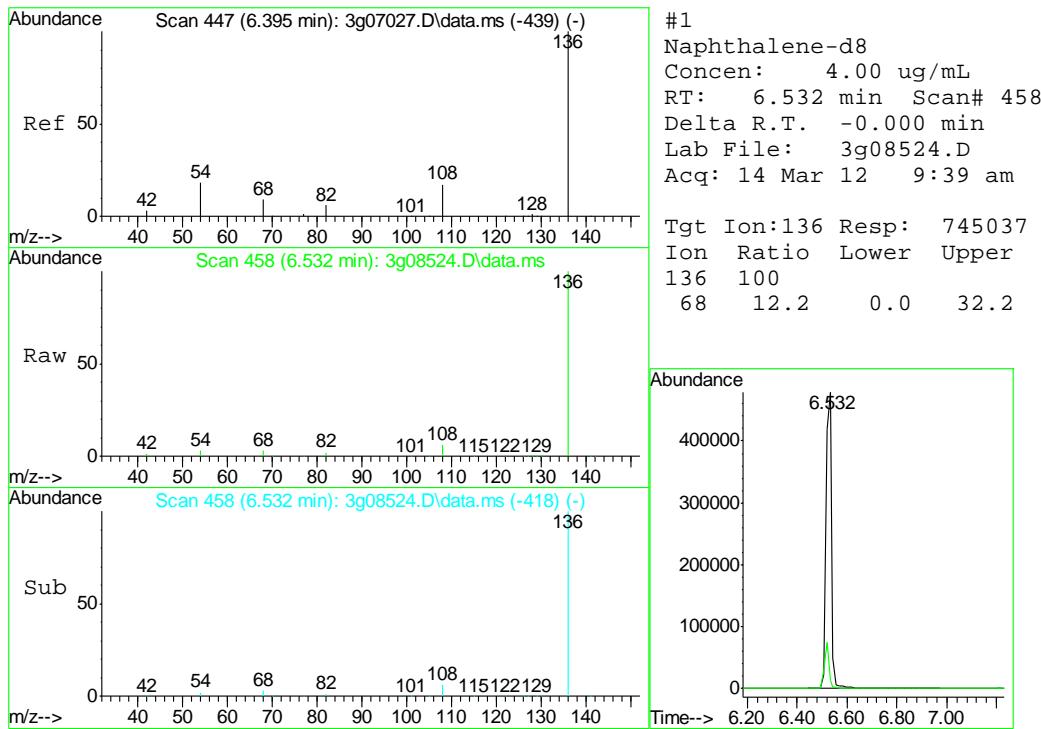
(#) = qualifier out of range (m) = manual integration (+) = signals summed

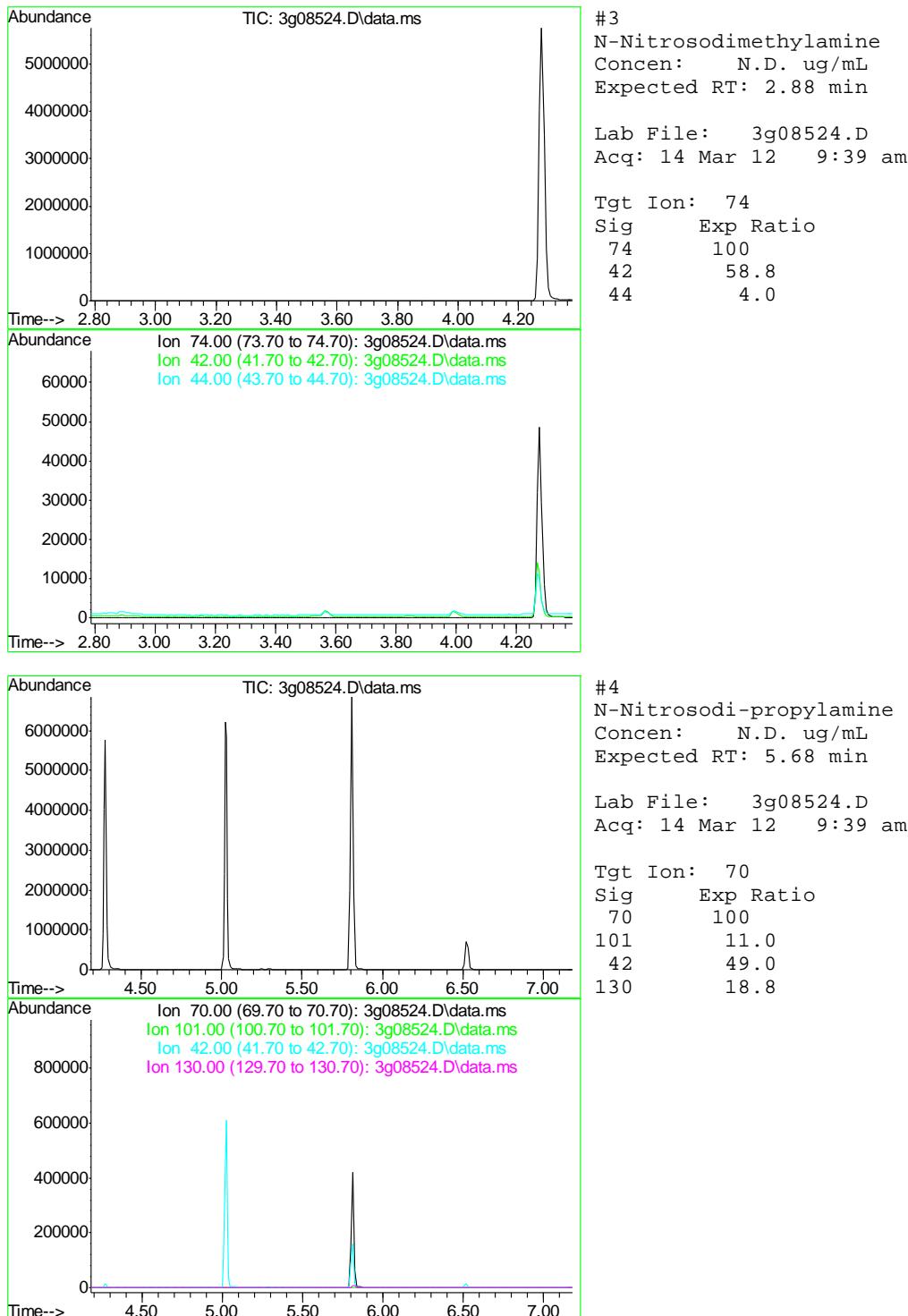
## Quantitation Report (QT Reviewed)

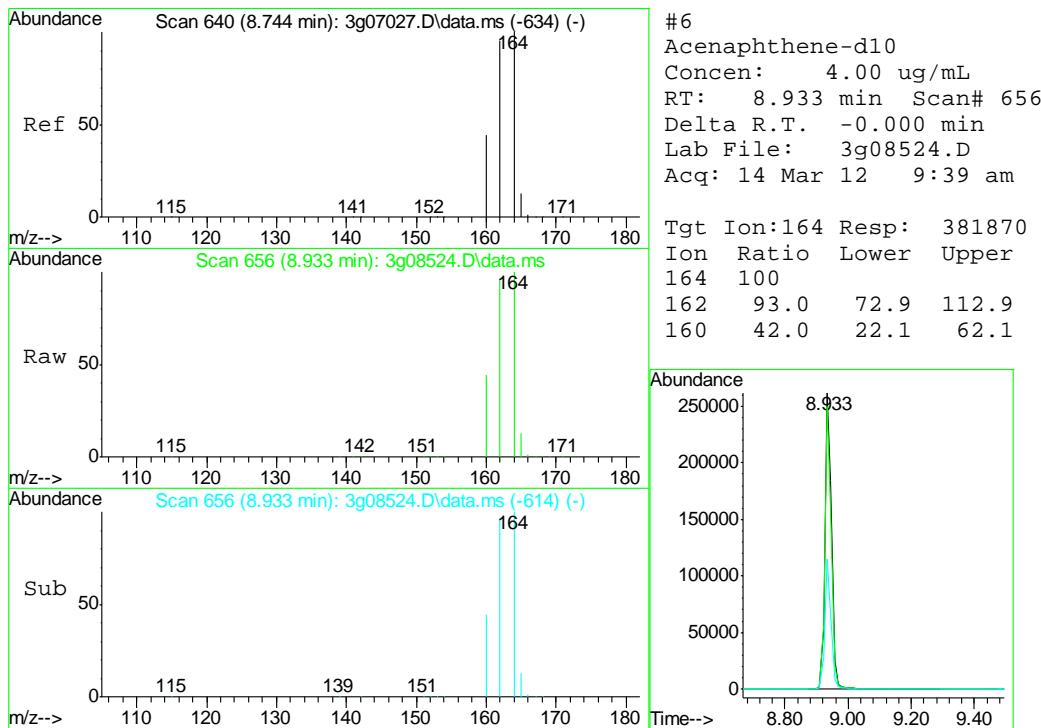
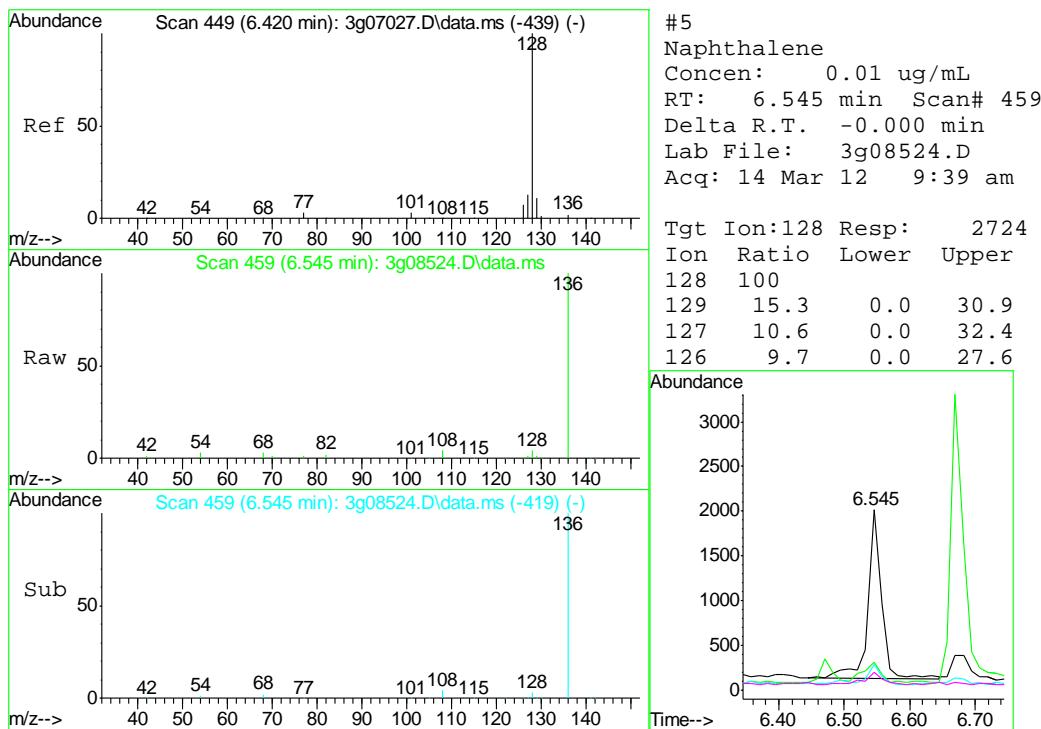
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 Data File : 3g08524.D  
 Acq On : 14 Mar 2012 9:39 am  
 Operator : DONC  
 Sample : OP5528-MB  
 Misc : OP5528,E3G348,30.00,,,1,1  
 ALS Vial : 4 Sample Multiplier: 1

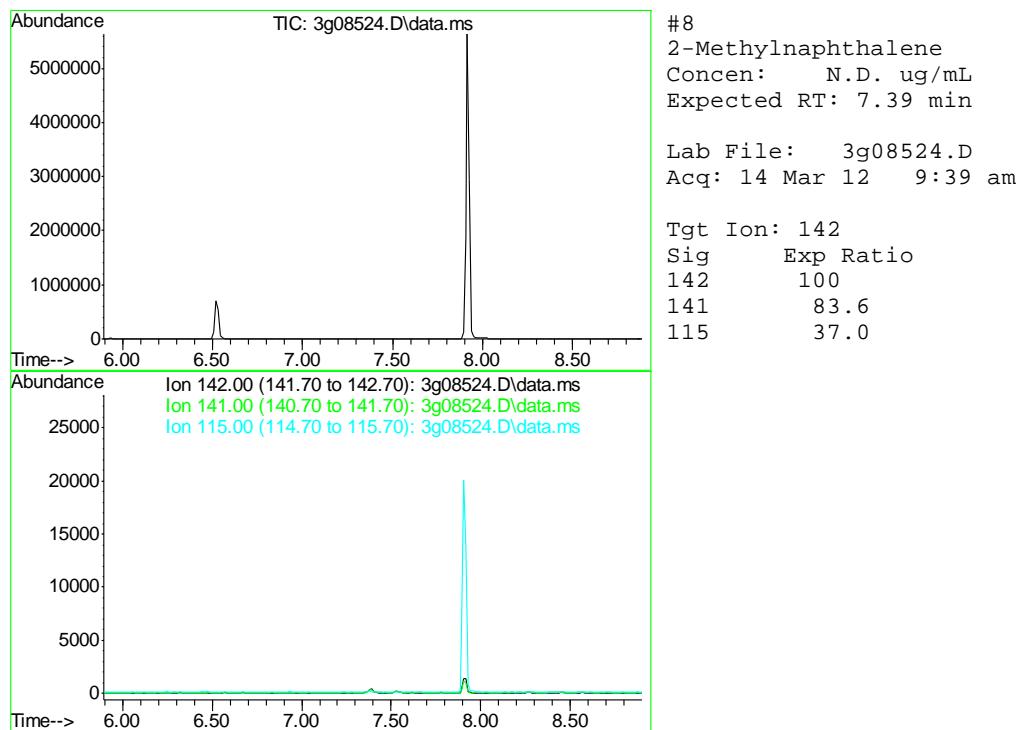
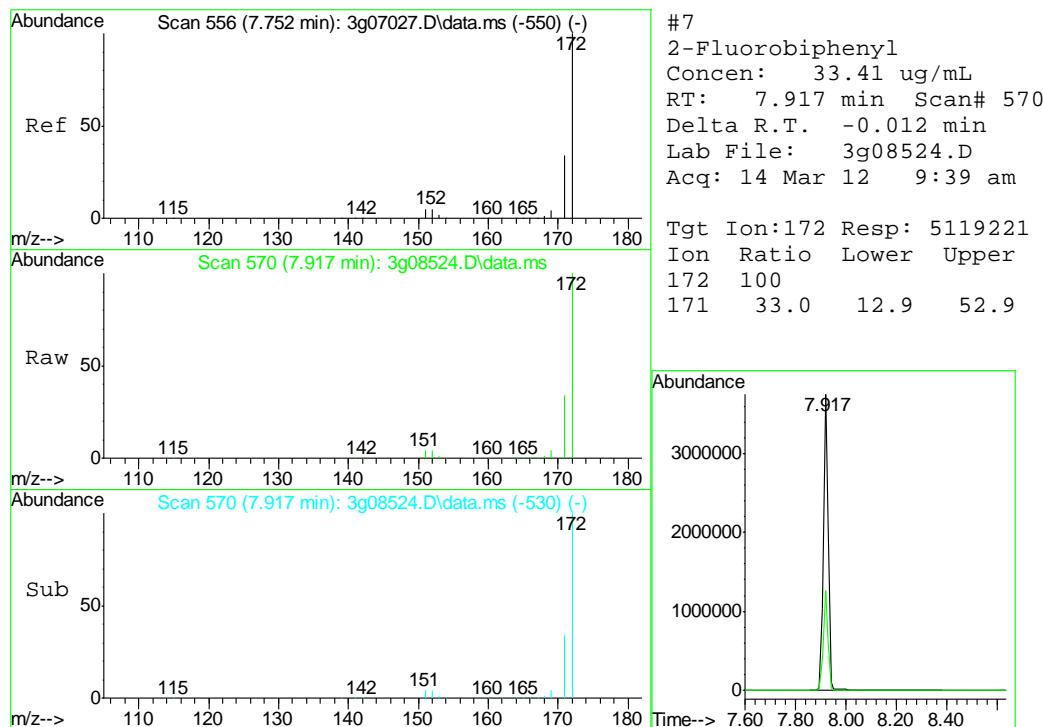
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 Quant Title : PAHSIM BASE  
 QLast Update : Mon Mar 12 09:19:25 2012  
 Response via : Initial Calibration

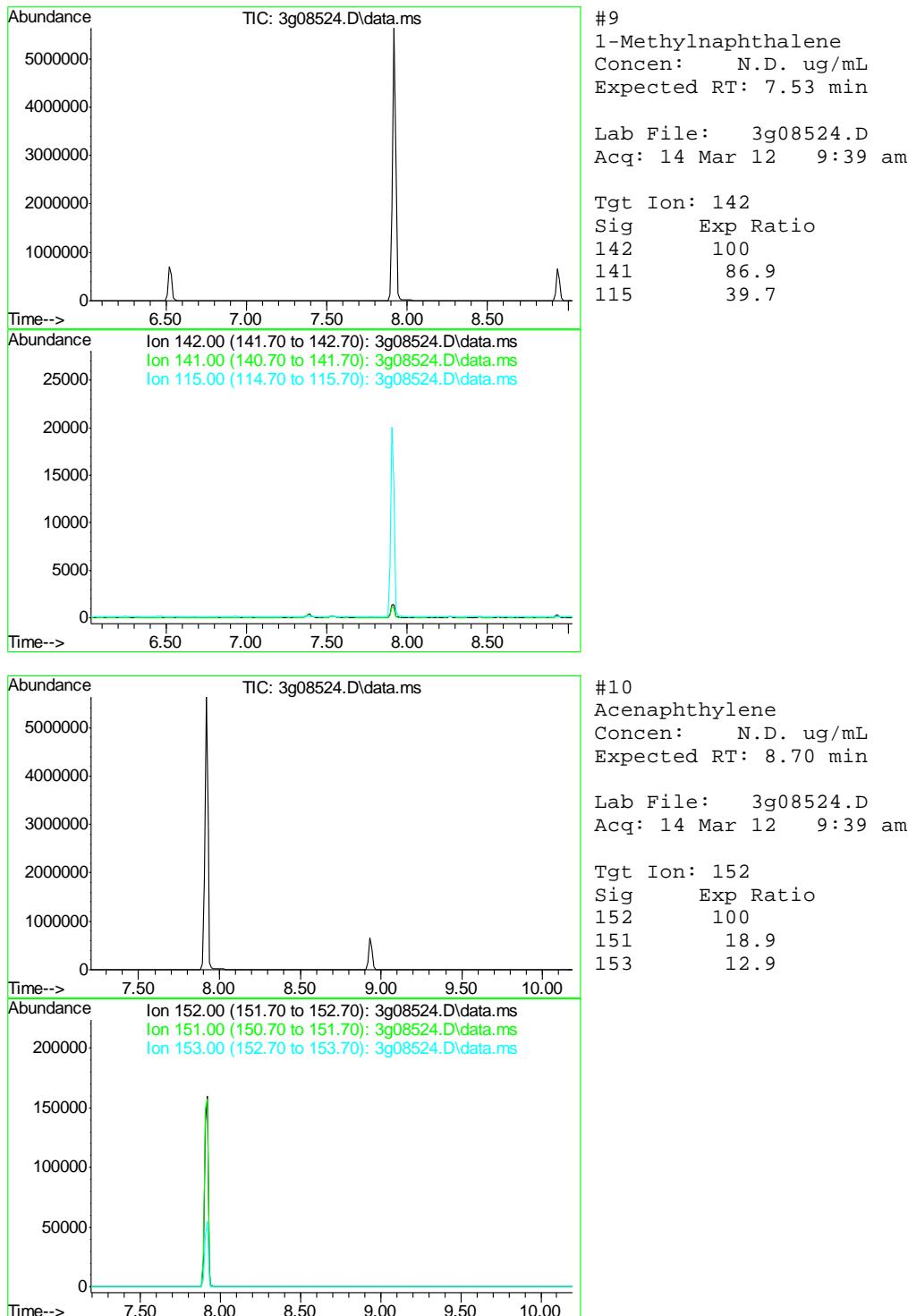


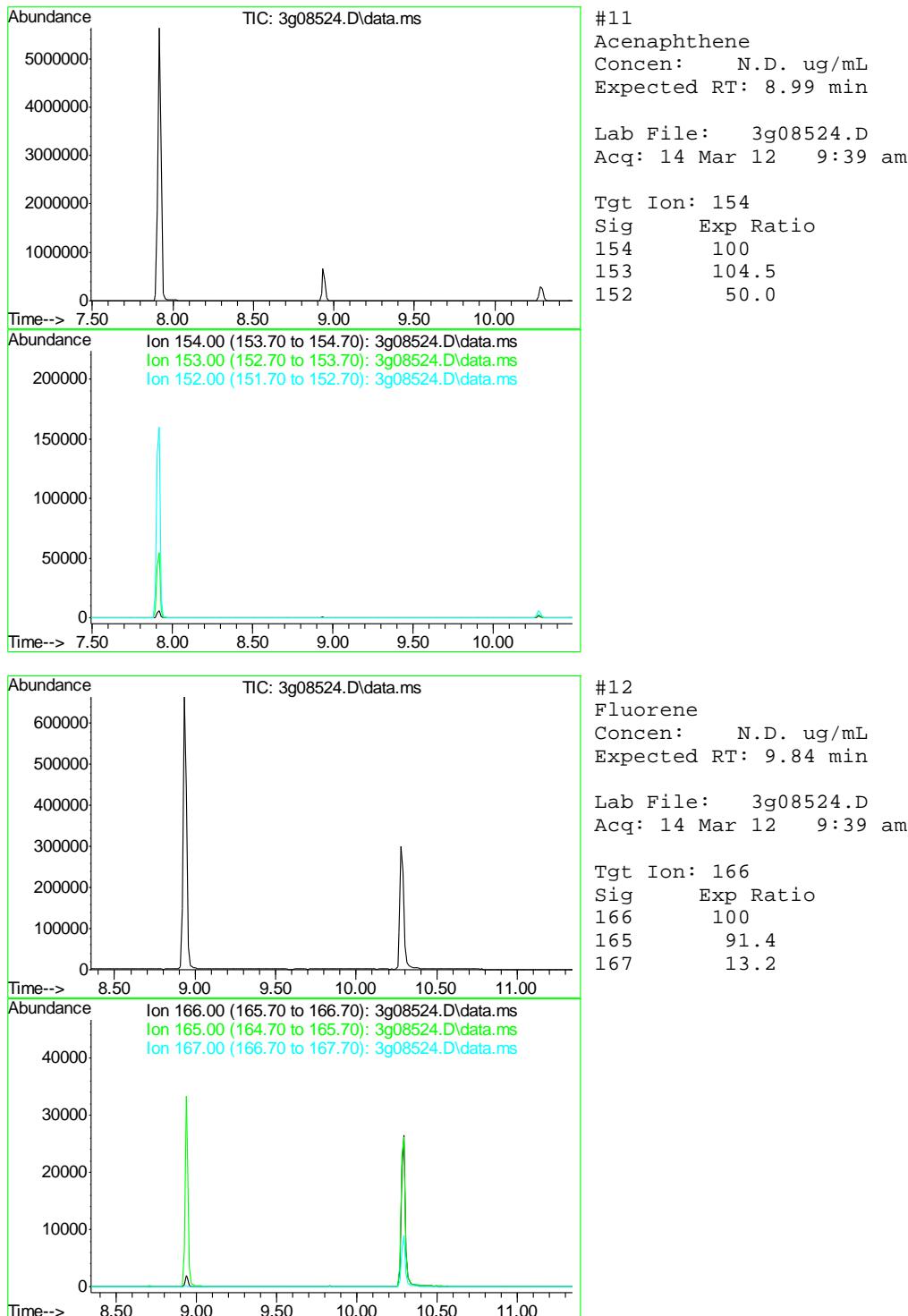


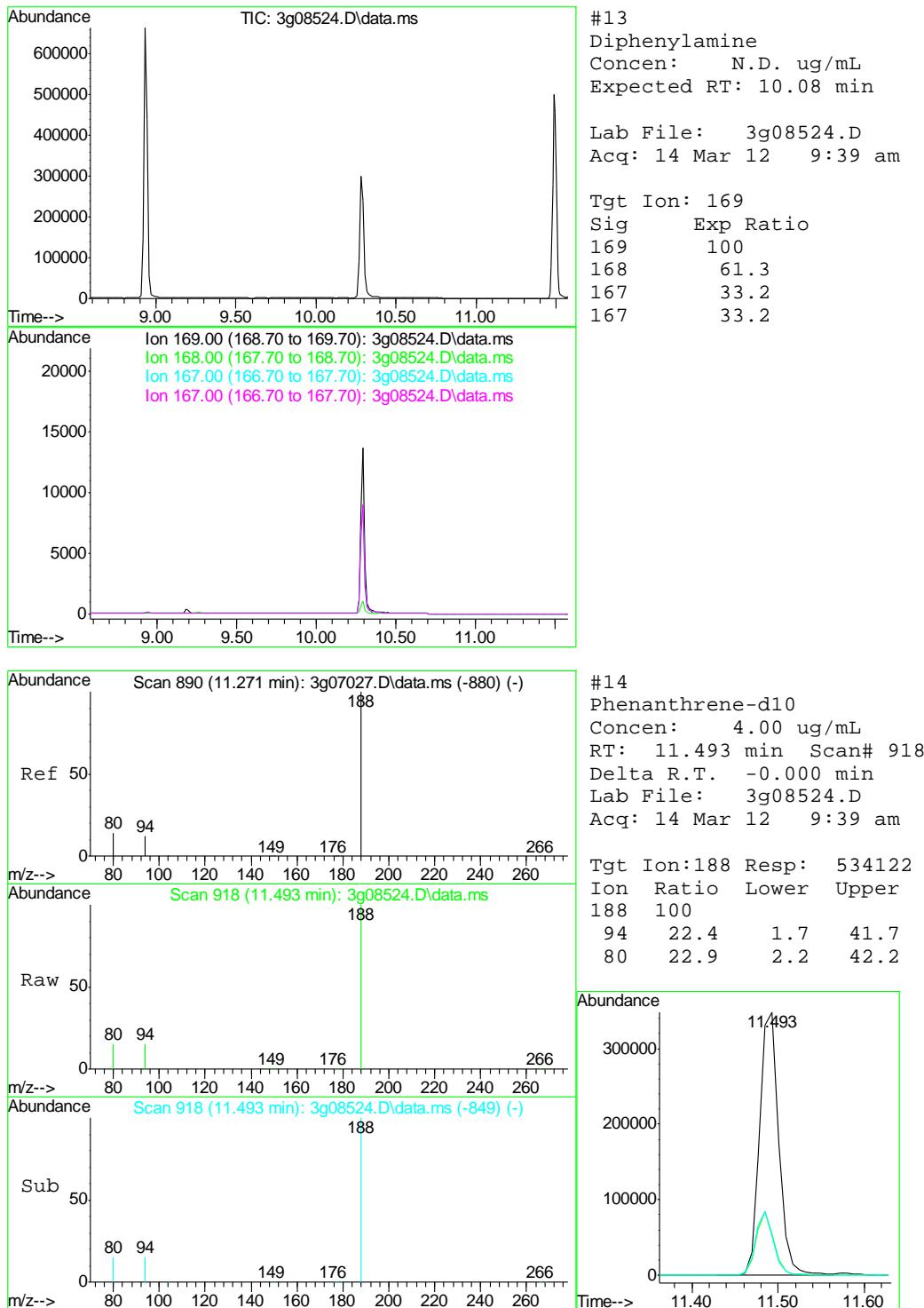


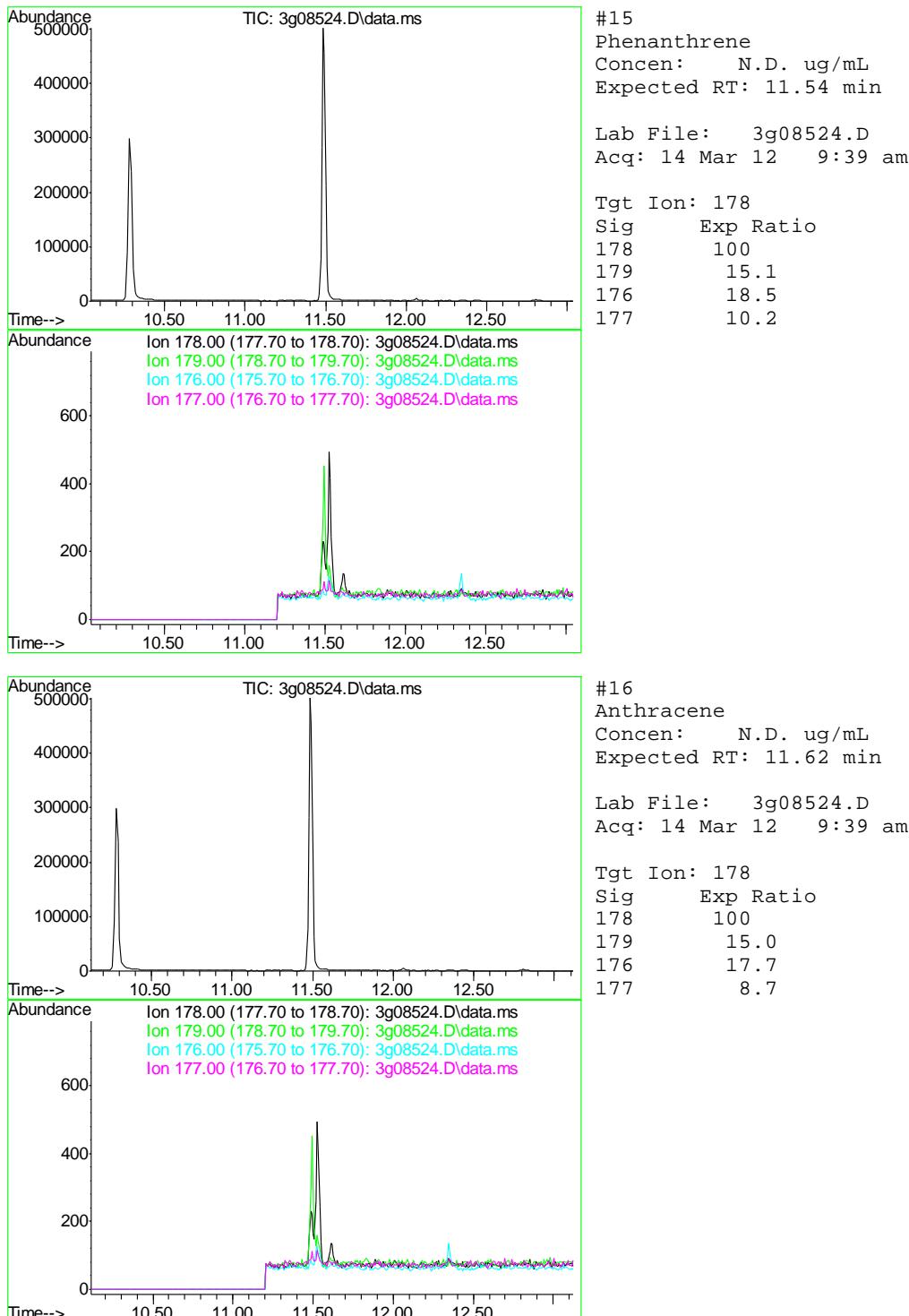


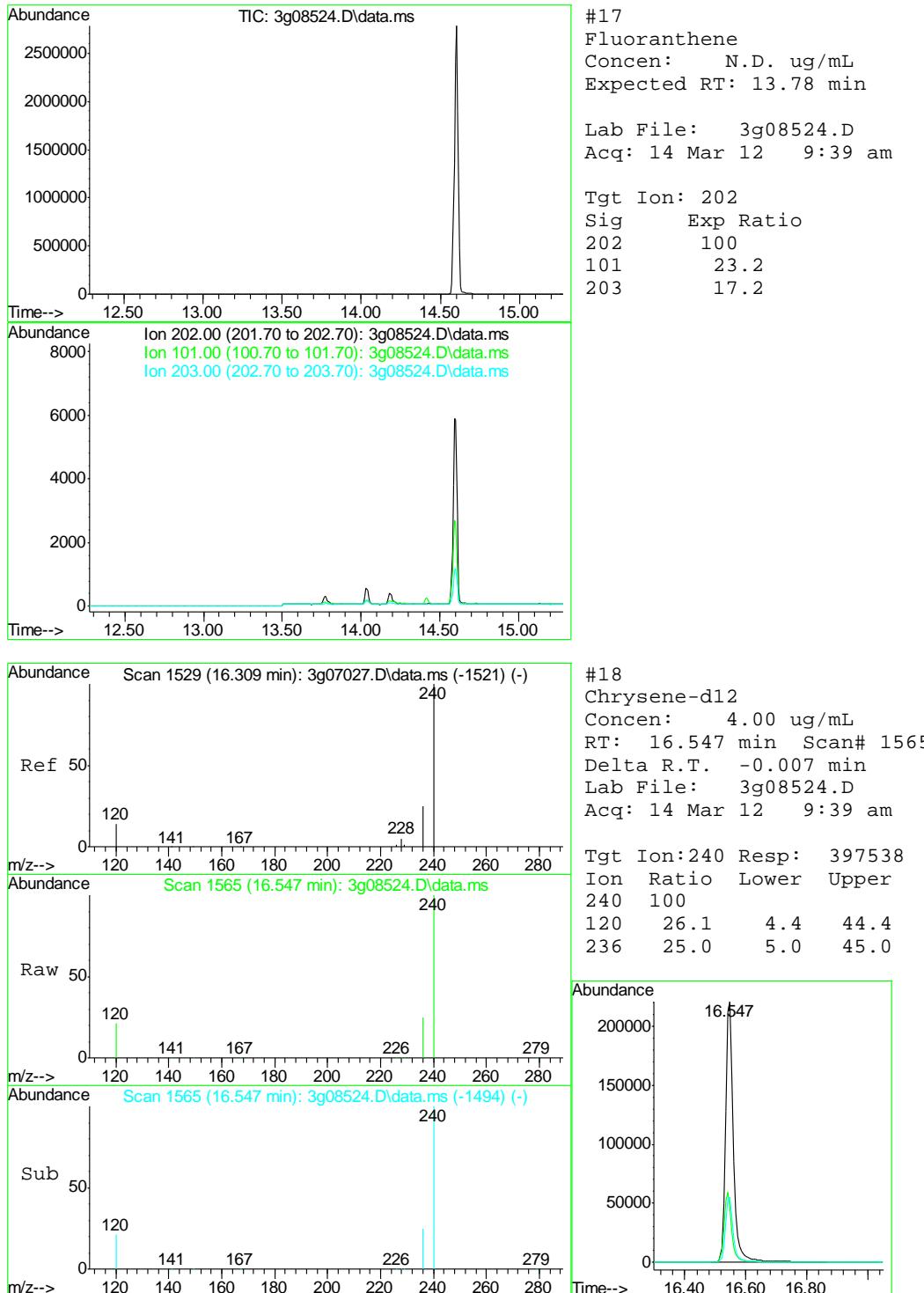


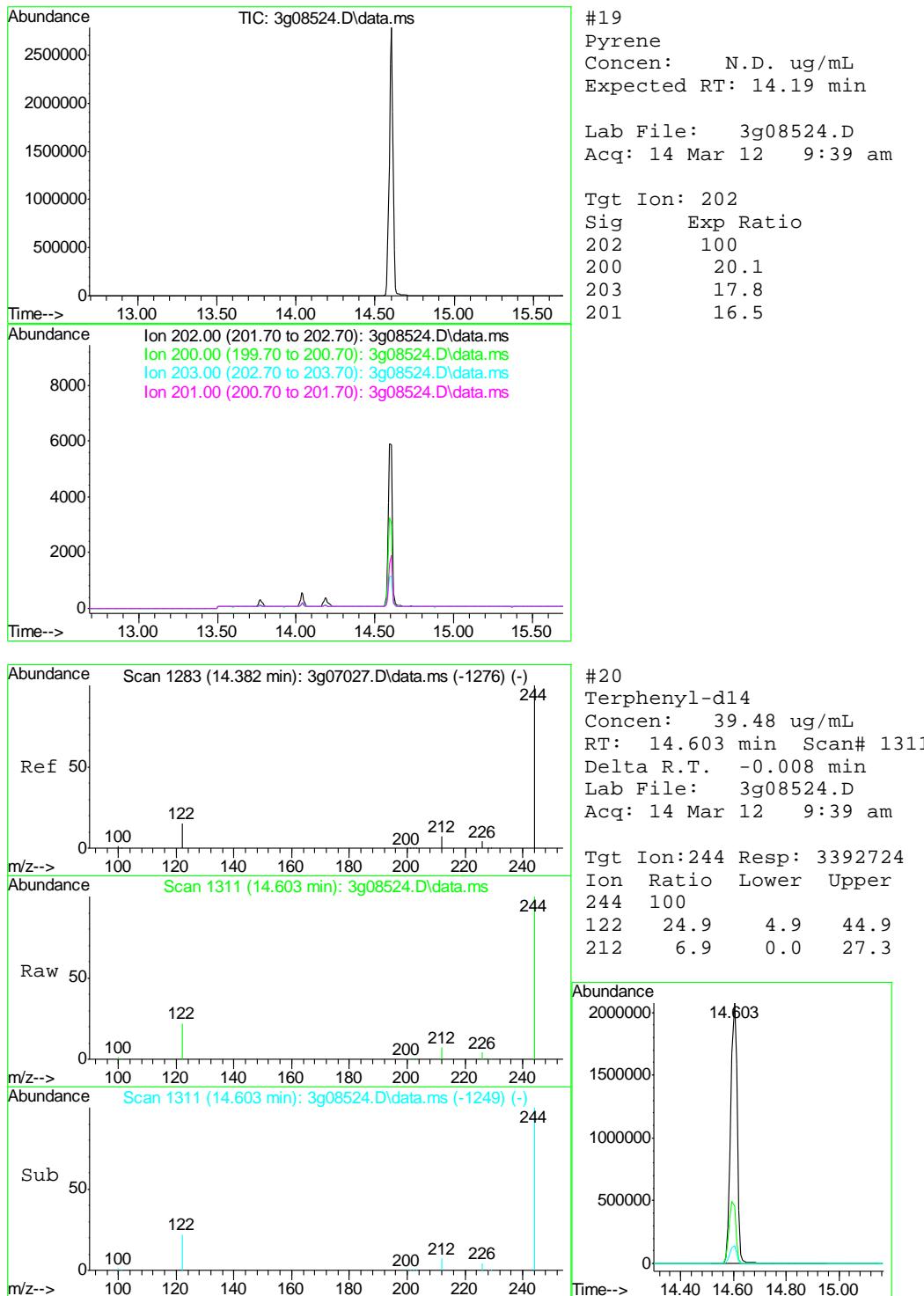


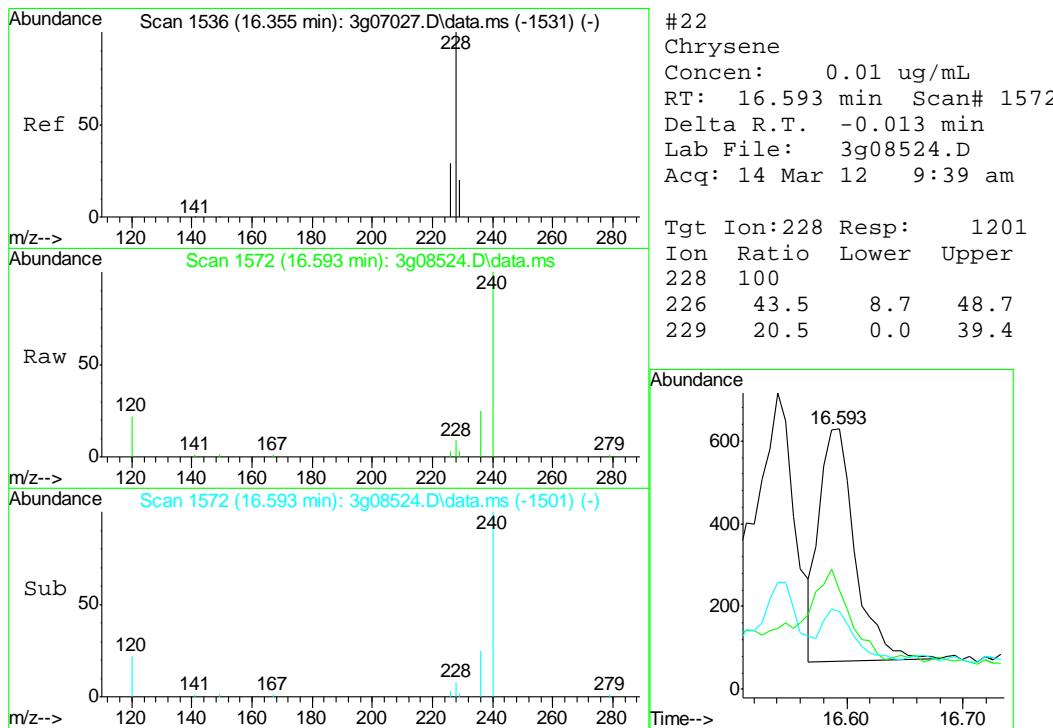
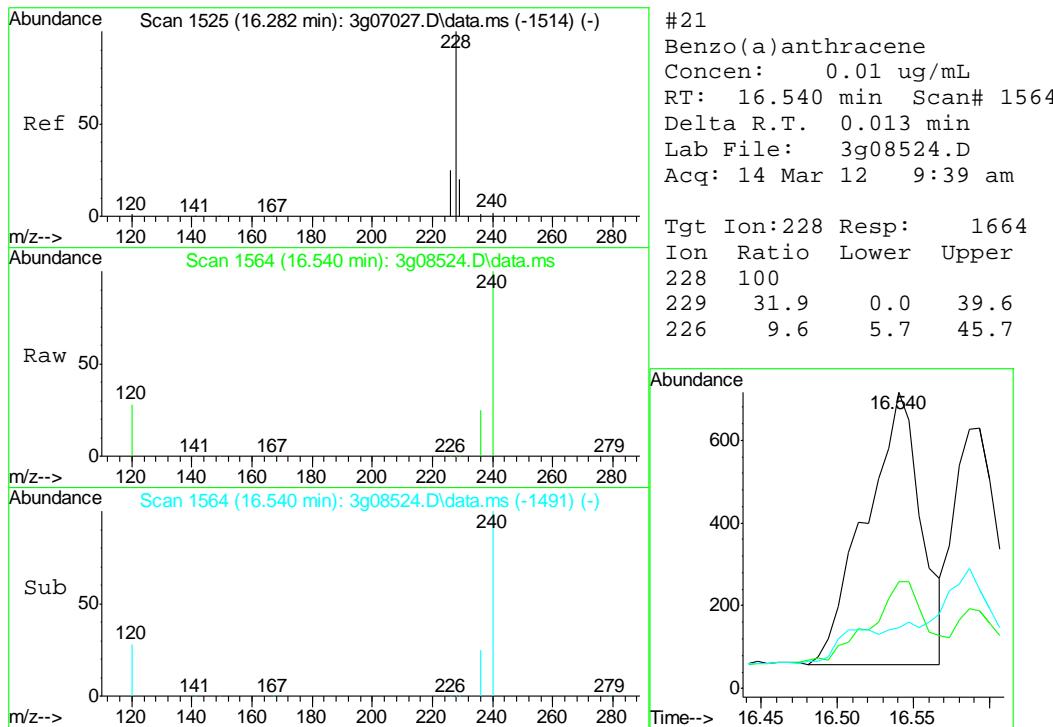


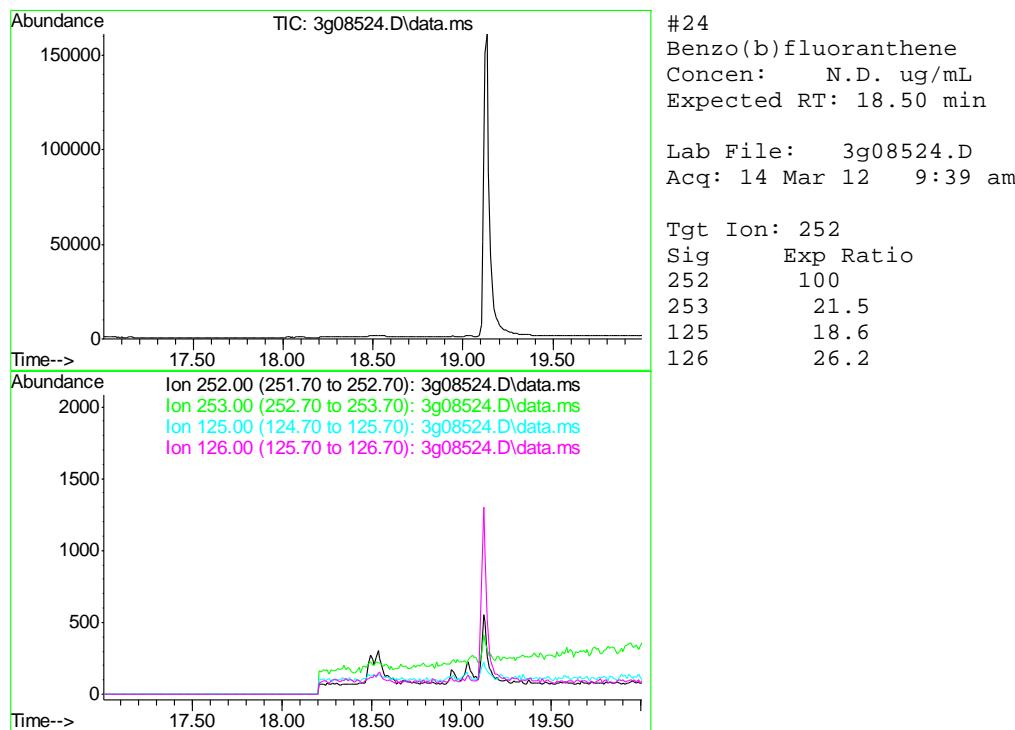
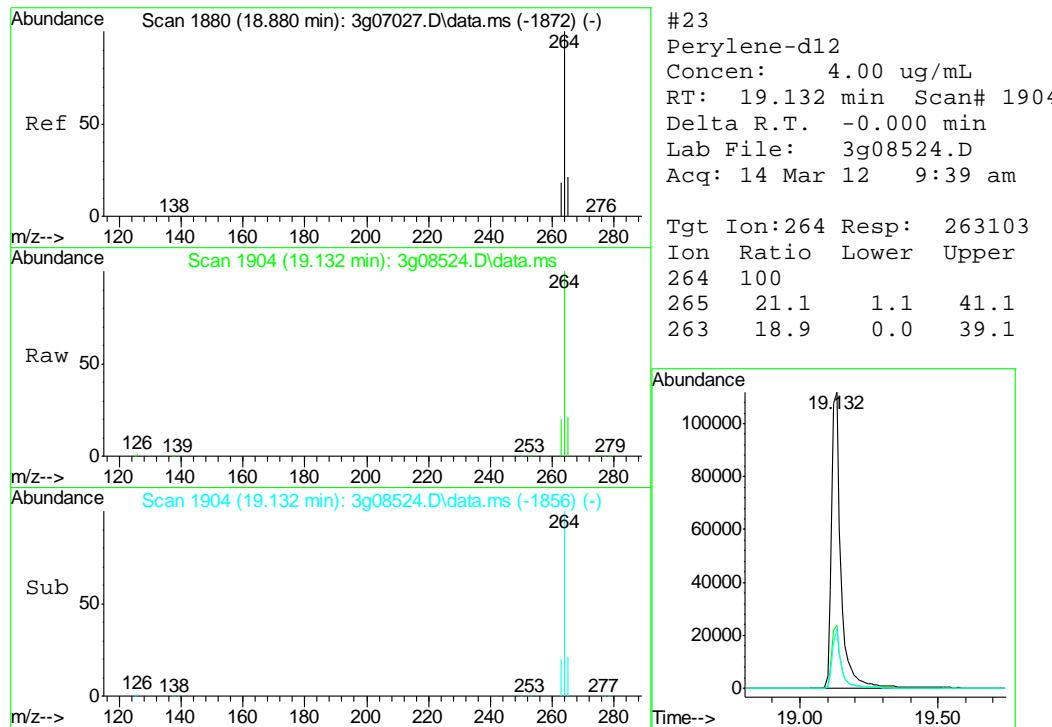


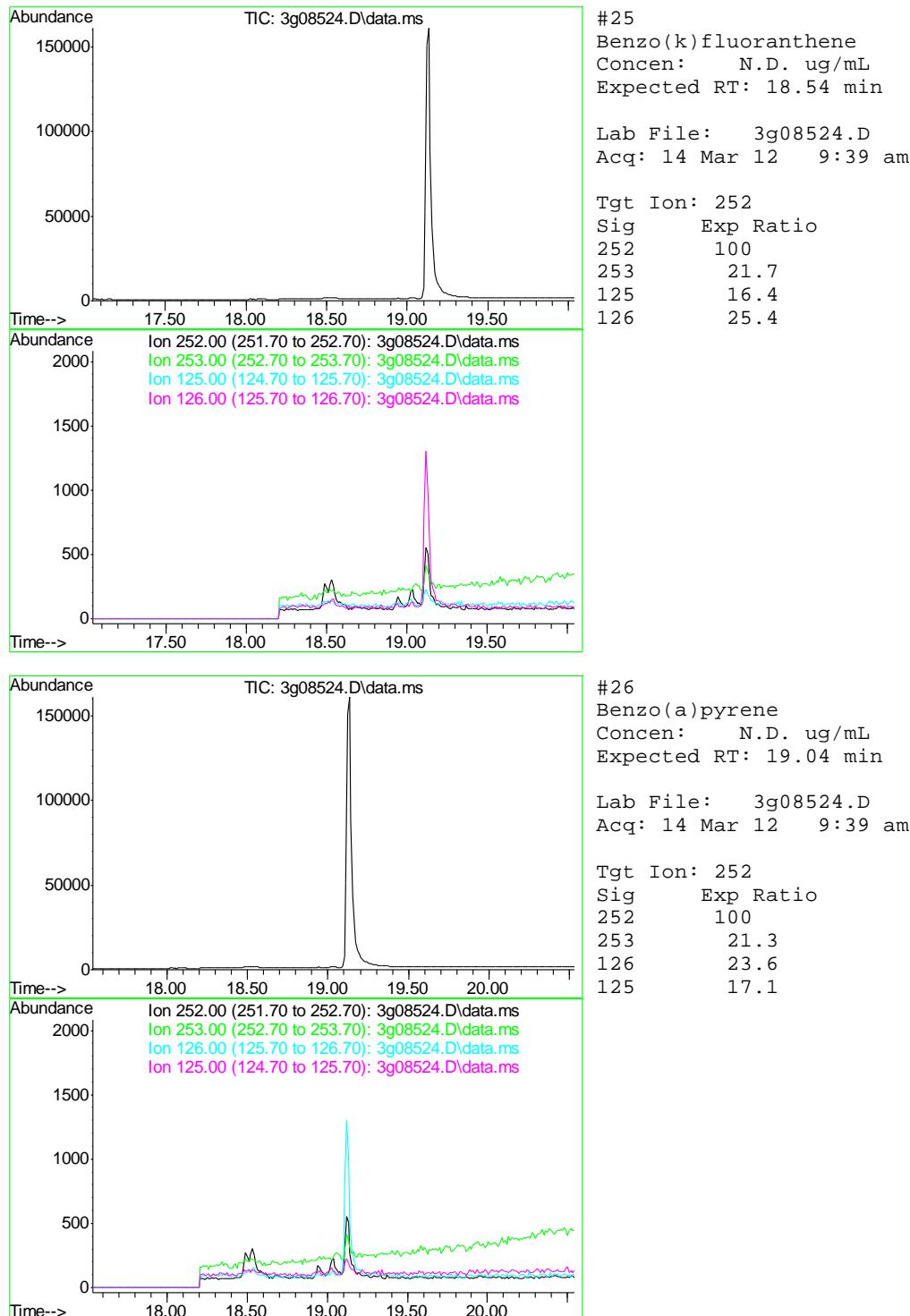


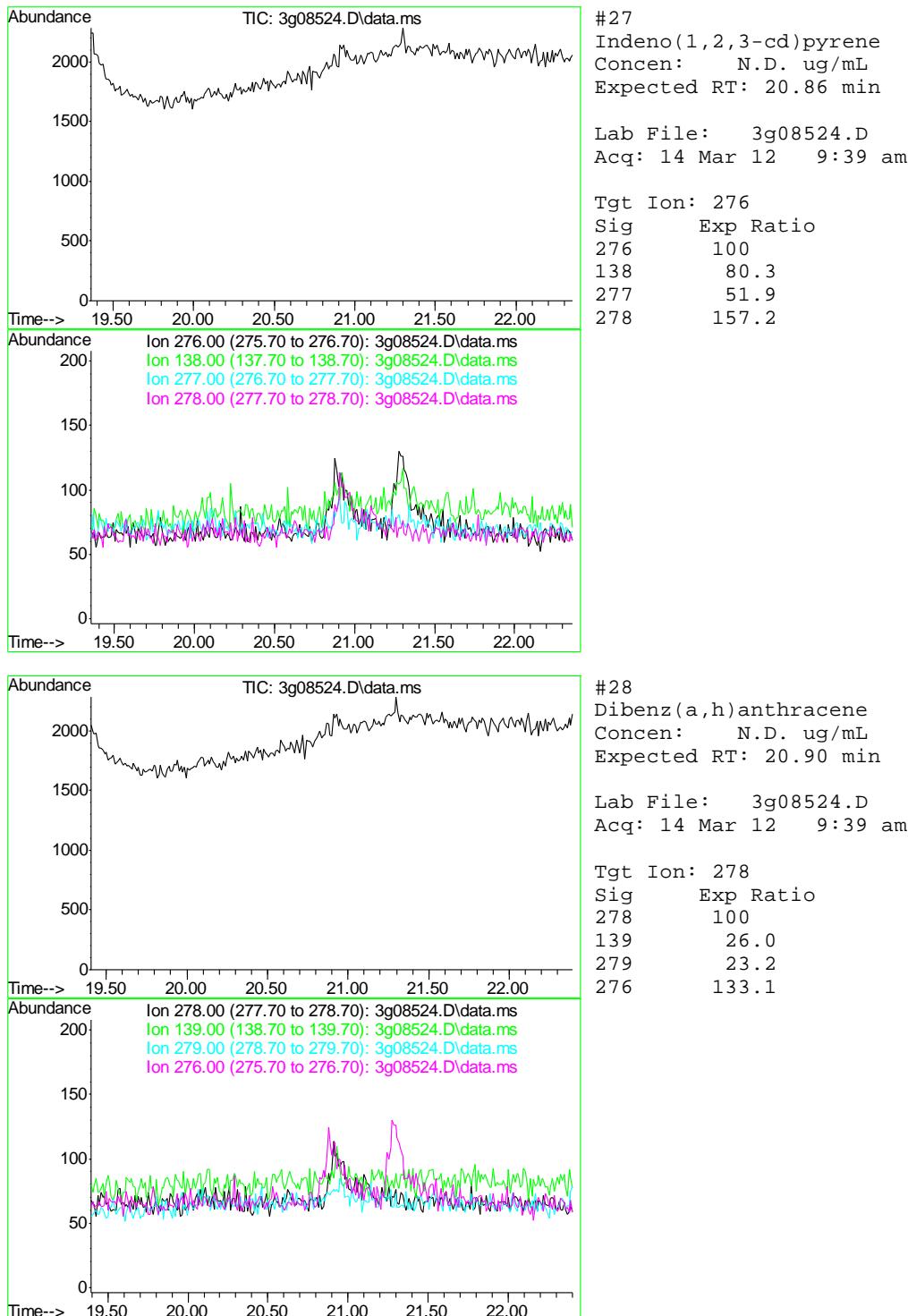


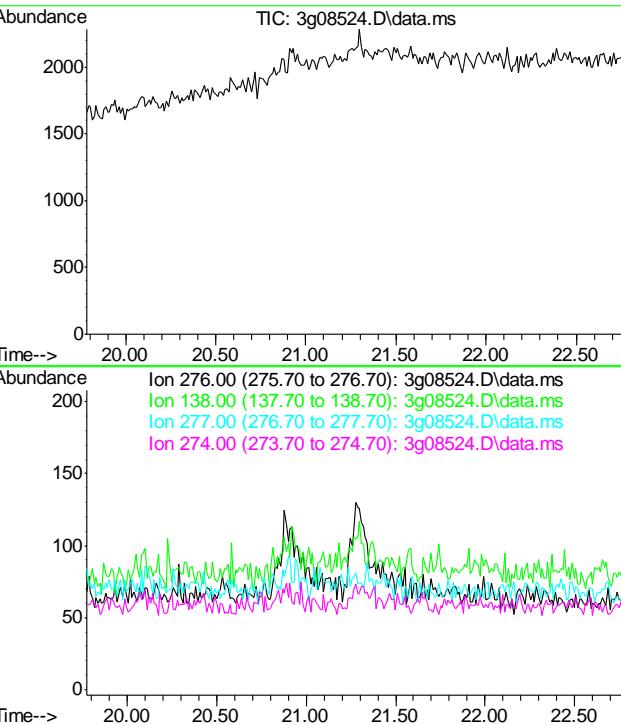












#29  
Benzo(g,h,i)perylene  
Concen: N.D. ug/mL  
Expected RT: 21.28 min

Lab File: 3g08524.D  
Acq: 14 Mar 12 9:39 am

Tgt Ion:	276
Sig	Exp Ratio
276	100
138	32.2
277	23.3
274	20.9

Abundance

Ion 276.00 (275.70 to 276.70): 3g08524.D\data.ms  
Ion 138.00 (137.70 to 138.70): 3g08524.D\data.ms  
Ion 277.00 (276.70 to 277.70): 3g08524.D\data.ms  
Ion 274.00 (273.70 to 274.70): 3g08524.D\data.ms

Time--> 20.00 20.50 21.00 21.50 22.00 22.50



## GC Volatiles

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### QC Data Summaries

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6

Includes the following where applicable:

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

**Method Blank Summary**

**Job Number:** D32609  
**Account:** XTOKWR XTO Energy  
**Project:** FRU 297-32A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
GGB858-MB	GB15295.D	1	03/13/12	SK	n/a	n/a	GGB858

The QC reported here applies to the following samples:

**Method:** SW846 8015B

D32609-1

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	109%      60-140%

9.1.1

9

## Blank Spike Summary

Page 1 of 1

Job Number: D32609

Account: XTOKWR XTO Energy

Project: FRU 297-32A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
GGB858-BS	GB15296.D	1	03/13/12	SK	n/a	n/a	GGB858

The QC reported here applies to the following samples:

Method: SW846 8015B

D32609-1

CAS No.	Compound	Spike mg/kg	BSP mg/kg	BSP %	Limits
	TPH-GRO (C6-C10)	110	112	102	70-130

CAS No.	Surrogate Recoveries	BSP	Limits
120-82-1	1,2,4-Trichlorobenzene	117%	60-140%

9.2.1

9

# Matrix Spike/Matrix Spike Duplicate Summary

Page 1 of 1

Job Number: D32609

Account: XTOKWR XTO Energy

Project: FRU 297-32A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
D32609-1MS	GB15298.D	1	03/13/12	SK	n/a	n/a	GGB858
D32609-1MSD	GB15299.D	1	03/13/12	SK	n/a	n/a	GGB858
D32609-1	GB15297.D	1	03/13/12	SK	n/a	n/a	GGB858

The QC reported here applies to the following samples:

Method: SW846 8015B

D32609-1

CAS No.	Compound	D32609-1		Spike	MS	MS	MSD	MSD	RPD	Limits Rec/RPD
		mg/kg	Q	mg/kg	mg/kg	%	mg/kg	%		
	TPH-GRO (C6-C10)	8.20	J	122	122	93	128	98	5	70-130/30

CAS No.	Surrogate Recoveries	MS	MSD	D32609-1	Limits
120-82-1	1,2,4-Trichlorobenzene	108%	111%	107%	60-140%

9.3.1

9



## GC Volatiles

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Raw Data

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Judy Nelson  
 03/14/12 09:45

## Quantitation Report (QT Reviewed)

Signal #1 : Y:\1\DATA\031312\GB15297.D\FID1A.CH Vial: 5  
 Signal #2 : Y:\1\DATA\031312\GB15297.D\FID2B.CH  
 Acq On : 13 Mar 2012 6:42 pm Operator: StephK  
 Sample : D32609-1, 50X Inst : GC/MS Ins  
 Misc : GC2671,GGB858,5.038,,100,5,1 Multiplr: 1.00  
 IntFile Signal #1: TVH1.E IntFile Signal #2: FB2.E  
 Quant Time: Mar 14 08:40:28 2012 Quant Results File: TB851GB851SOIL.RES

Quant Method : C:\MSDCHEM\1...\TB851GB851SOIL.M (Chemstation Integrator)  
 Title : 8015B/8021B TVH/BTEX  
 Last Update : Wed Mar 14 08:40:04 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
----------	------	----------	------	-------

## System Monitoring Compounds

2) S	1,2,4-Trichlorobenzene	14.34	3221074	106.774 %	m
10) S	1,2,4-Trichlorobenzene (P)	14.34	25943932	111.481 %	

## Target Compounds

1) H	TVH-Gasoline	7.26	10666507	0.148 mg/L
4) T	Methyl-t-butyl-ether	0.00	0	N.D. ug/L d
5) T	Benzene	4.09	504716	0.912 ug/L
6) T	Toluene	7.60	1338355	2.445 ug/L
7) T	Ethylbenzene	10.24	287853	0.629 ug/L
8) T	m,p-Xylene	10.42	1527112	2.727 ug/L
9) T	o-Xylene	10.94	295336	0.644 ug/L
11) T	Naphthalene	14.52	1455364	5.570 ug/L

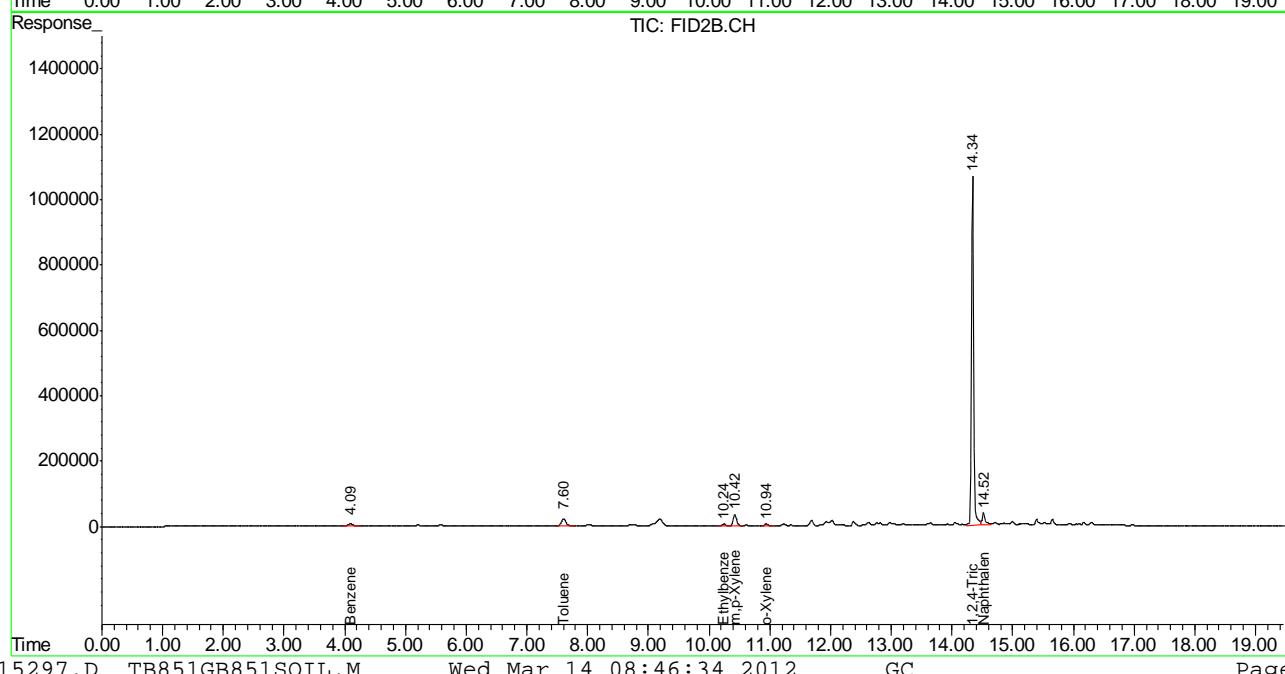
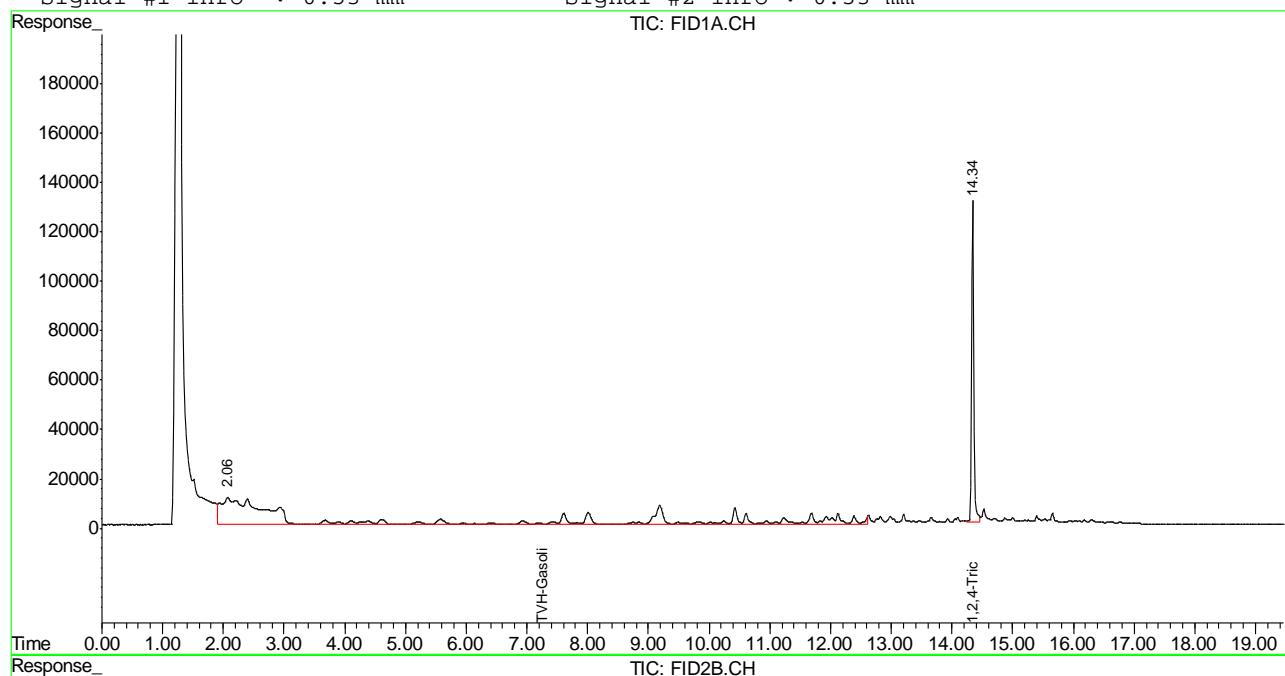
(f)=RT Delta > 1/2 Window (m)=manual int.  
 GB15297.D TB851GB851SOIL.M Wed Mar 14 08:46:34 2012 GC

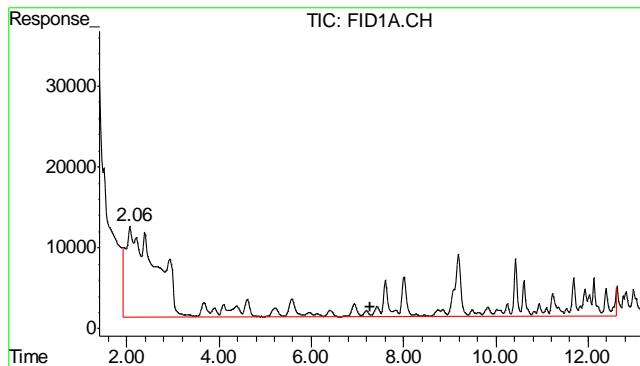
## Quantitation Report (QT Reviewed)

Signal #1 : Y:\1\DATA\031312\GB15297.D\FID1A.CH Vial: 5  
 Signal #2 : Y:\1\DATA\031312\GB15297.D\FID2B.CH  
 Acq On : 13 Mar 2012 6:42 pm Operator: StephK  
 Sample : D32609-1, 50X Inst : GC/MS Ins  
 Misc : GC2671,GGB858,5.038,,100,5,1 Multiplr: 1.00  
 IntFile Signal #1: TVH1.E IntFile Signal #2: FB2.E  
 Quant Time: Mar 14 8:43 2012 Quant Results File: TB851GB851SOIL.RES

Quant Method : C:\MSDCHEM\1...\TB851GB851SOIL.M (Chemstation Integrator)  
 Title : 8015B/8021B TVH/BTEX  
 Last Update : Wed Mar 14 08:40:04 2012  
 Response via : Multiple Level Calibration  
 DataAcq Meth : TVB4.M

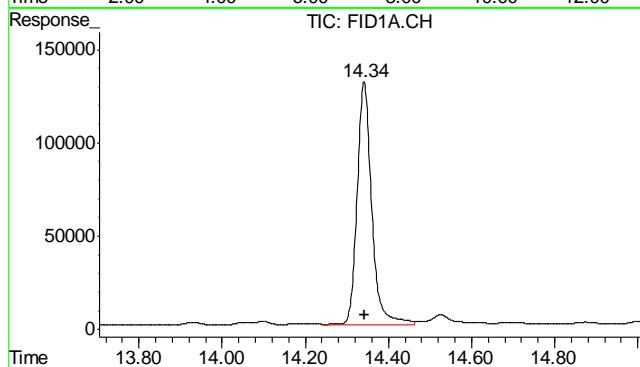
Volume Inj. :  
 Signal #1 Phase : DB-624 Signal #2 Phase: DB-624  
 Signal #1 Info : 0.53 mm Signal #2 Info : 0.53 mm





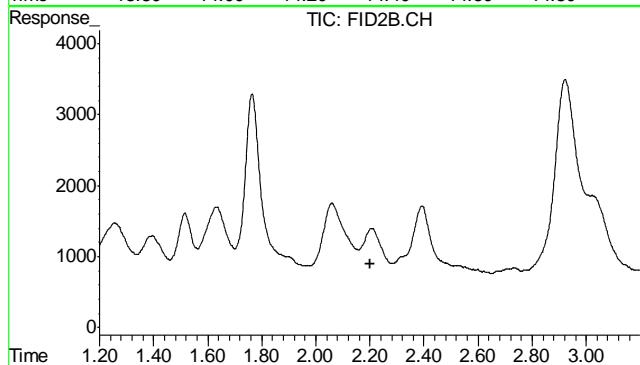
#1 TVH-Gasoline

R.T.: 7.265 min  
Delta R.T.: 0.000 min  
Response: 10666507  
Conc: 0.15 mg/L m



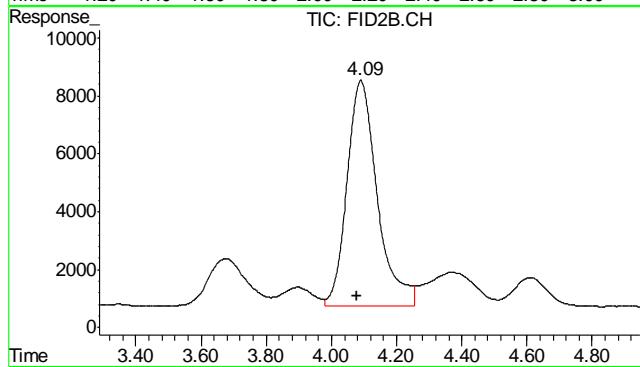
#2 1,2,4-Trichlorobenzene

R.T.: 14.341 min  
Delta R.T.: 0.000 min  
Response: 3221074  
Conc: 106.77 % m



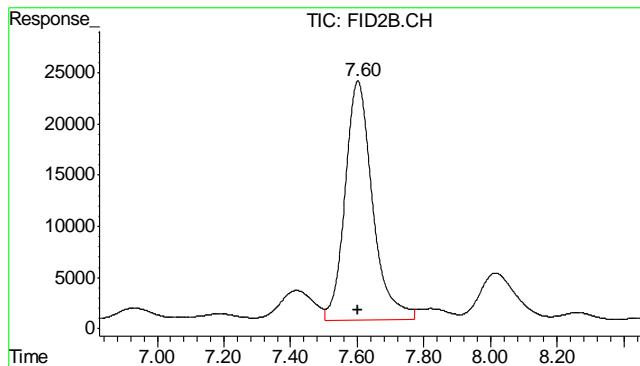
#4 Methyl-t-butyl-ether

R.T.: 0.000 min  
Exp R.T.: 2.199 min  
Response: 0  
Conc: N.D.

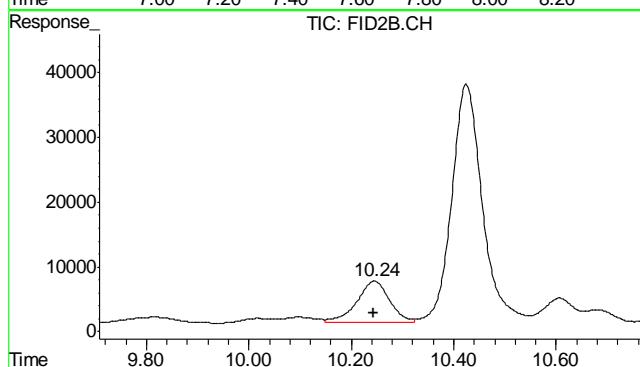


#5 Benzene

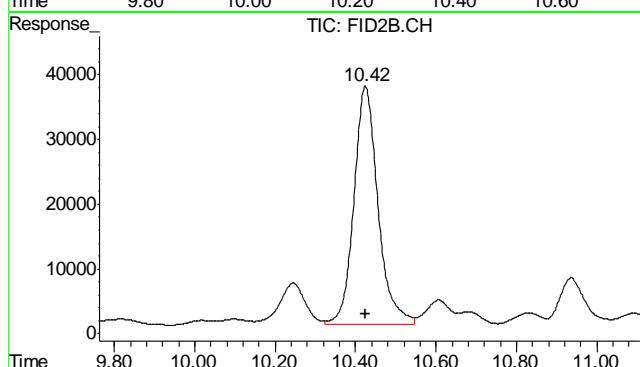
R.T.: 4.090 min  
Delta R.T.: 0.012 min  
Response: 504716  
Conc: 0.91 ug/L



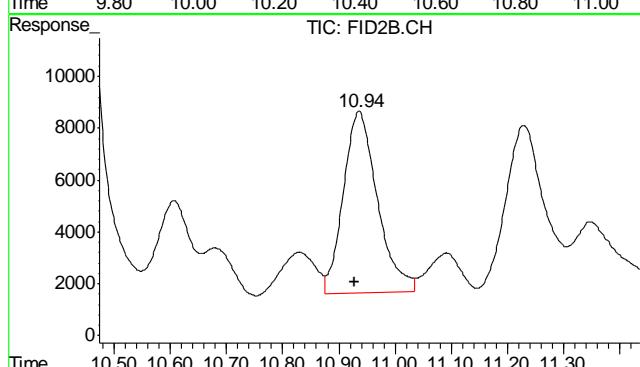
#6 Toluene  
R.T.: 7.603 min  
Delta R.T.: 0.002 min  
Response: 1338355  
Conc: 2.44 ug/L



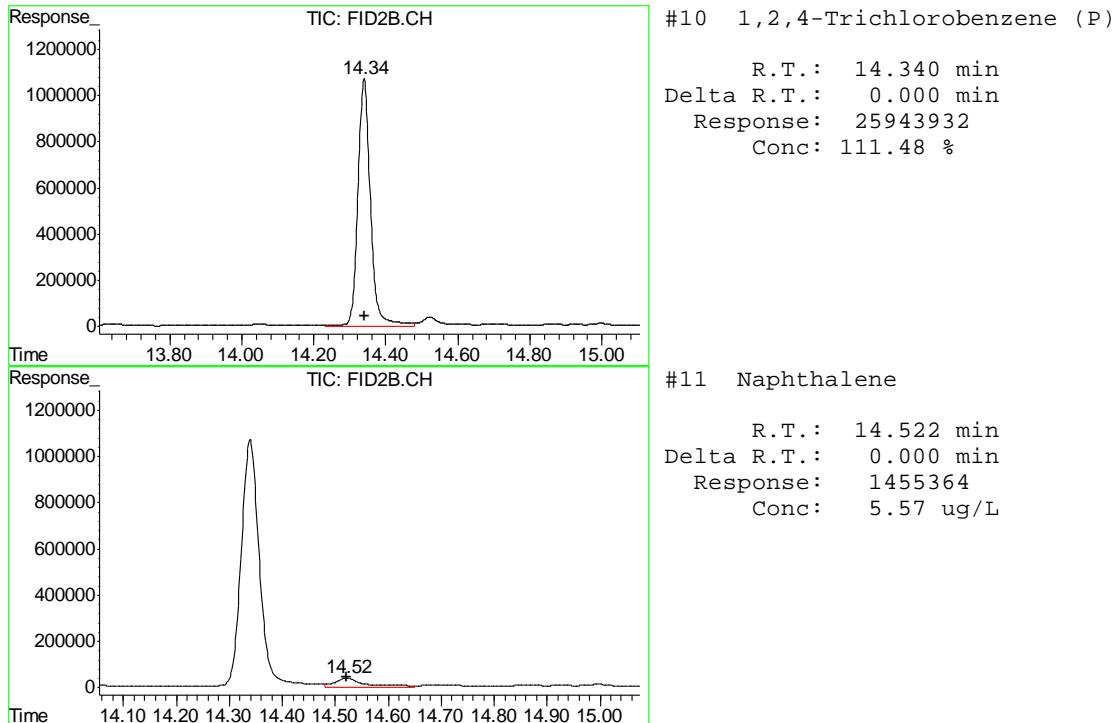
#7 Ethylbenzene  
R.T.: 10.245 min  
Delta R.T.: 0.000 min  
Response: 287853  
Conc: 0.63 ug/L



#8 m,p-Xylene  
R.T.: 10.425 min  
Delta R.T.: -0.002 min  
Response: 1527112  
Conc: 2.73 ug/L



#9 o-Xylene  
R.T.: 10.936 min  
Delta R.T.: 0.009 min  
Response: 295336  
Conc: 0.64 ug/L



10.1.1

10

Judy Nelson  
 03/14/12 09:45

## Quantitation Report (QT Reviewed)

Signal #1 : Y:\1\DATA\031312\GB15295.D\FID1A.CH Vial: 3  
 Signal #2 : Y:\1\DATA\031312\GB15295.D\FID2B.CH  
 Acq On : 13 Mar 2012 5:30 pm Operator: StephK  
 Sample : MB Inst : GC/MS Ins  
 Misc : GC2671,GGB858,5.000,,100,5,1 Multiplr: 1.00  
 IntFile Signal #1: TVH1.E IntFile Signal #2: FB2.E  
 Quant Time: Mar 14 08:40:20 2012 Quant Results File: TB851GB851SOIL.RES

Quant Method : C:\MSDCHEM\1...\TB851GB851SOIL.M (Chemstation Integrator)  
 Title : 8015B/8021B TVH/BTEX  
 Last Update : Wed Mar 14 08:40:04 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	14.34	3279322	108.705 %	m
10) S	1,2,4-Trichlorobenzene (P)	14.34	26080780	112.069 %	

## Target Compounds

1) H	TVH-Gasoline	7.26	5696664	<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	7.60	218482	0.399	ug/L
7) T	Ethylbenzene	0.00	0	N.D.	ug/L d
8) T	m,p-Xylene	10.42	220790	0.394	ug/L
9) T	o-Xylene	10.93	88601	0.193	ug/L
11) T	Naphthalene	14.52	378695	1.449	ug/L

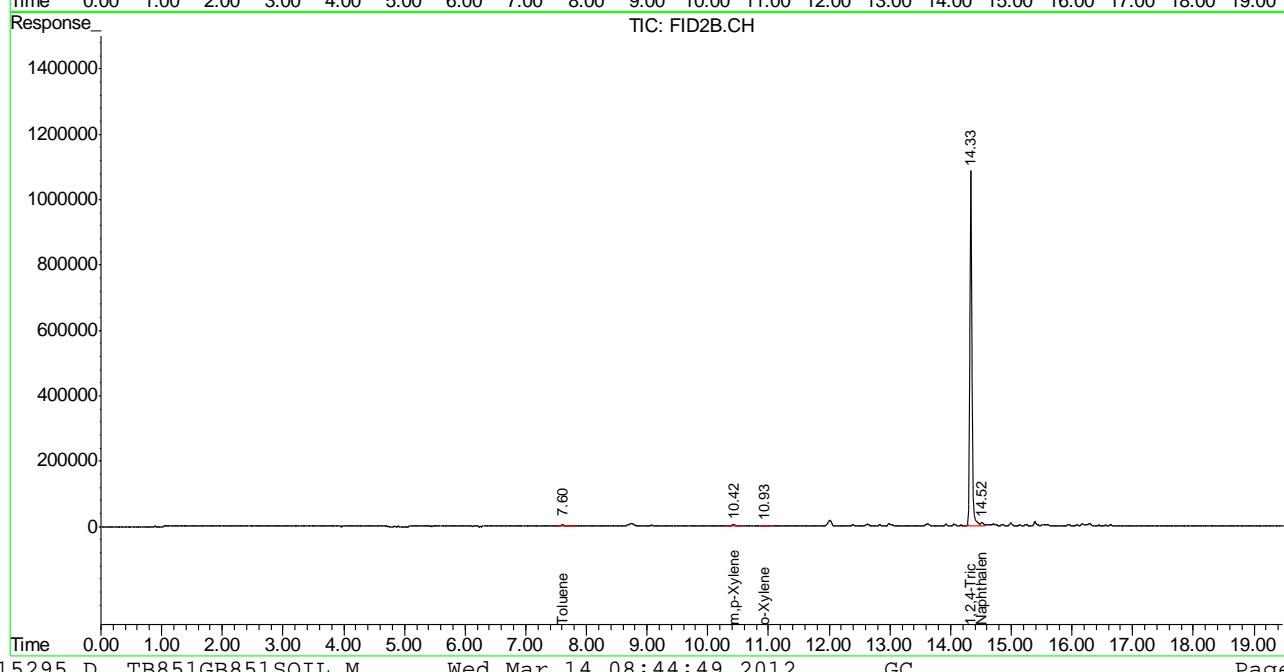
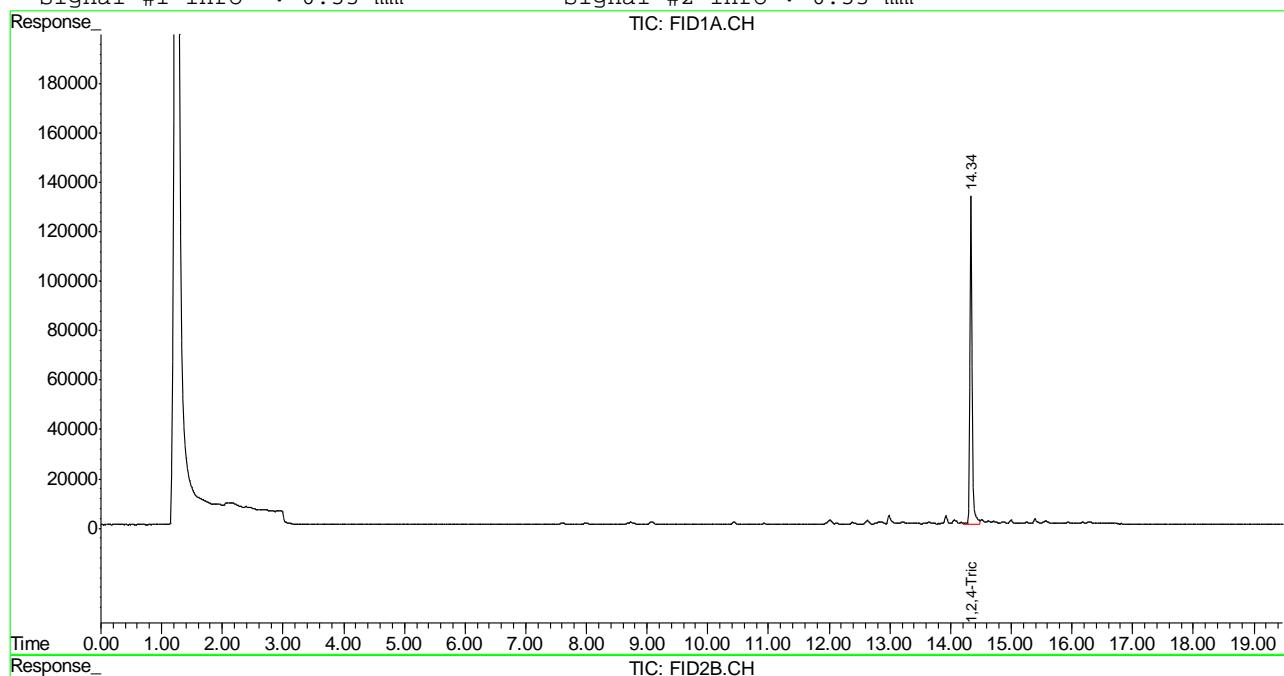
(f)=RT Delta > 1/2 Window (m)=manual int.  
 GB15295.D TB851GB851SOIL.M Wed Mar 14 08:44:49 2012 GC

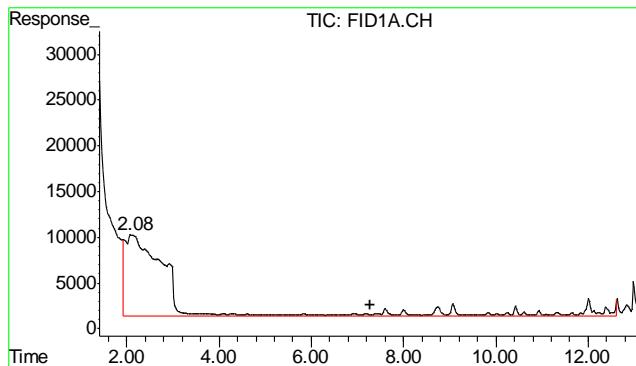
## Quantitation Report (QT Reviewed)

Signal #1 : Y:\1\DATA\031312\GB15295.D\FID1A.CH Vial: 3  
 Signal #2 : Y:\1\DATA\031312\GB15295.D\FID2B.CH  
 Acq On : 13 Mar 2012 5:30 pm Operator: StephK  
 Sample : MB Inst : GC/MS Ins  
 Misc : GC2671,GGB858,5.000,,100,5,1 Multiplr: 1.00  
 IntFile Signal #1: TVH1.E IntFile Signal #2: FB2.E  
 Quant Time: Mar 14 8:43 2012 Quant Results File: TB851GB851SOIL.RES

Quant Method : C:\MSDCHEM\1...\TB851GB851SOIL.M (Chemstation Integrator)  
 Title : 8015B/8021B TVH/BTEX  
 Last Update : Wed Mar 14 08:40:04 2012  
 Response via : Multiple Level Calibration  
 DataAcq Meth : TVB4.M

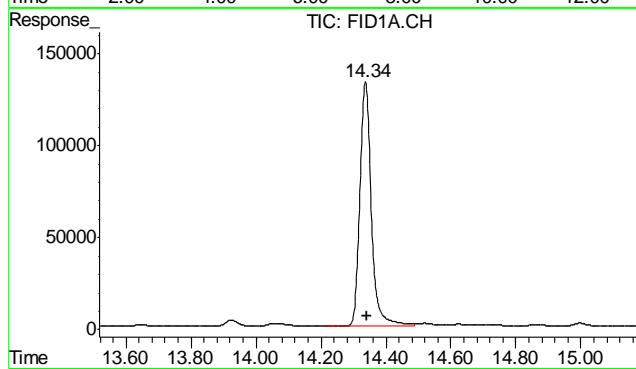
Volume Inj. :  
 Signal #1 Phase : DB-624 Signal #2 Phase: DB-624  
 Signal #1 Info : 0.53 mm Signal #2 Info : 0.53 mm





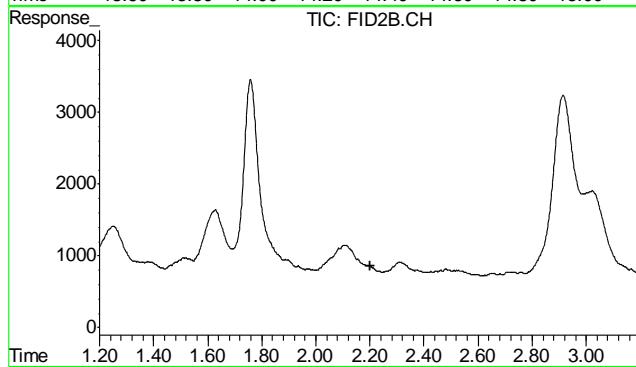
#1 TVH-Gasoline

R.T.: 7.265 min  
 Delta R.T.: 0.000 min  
 Response: 5696664  
 Conc: N.D.



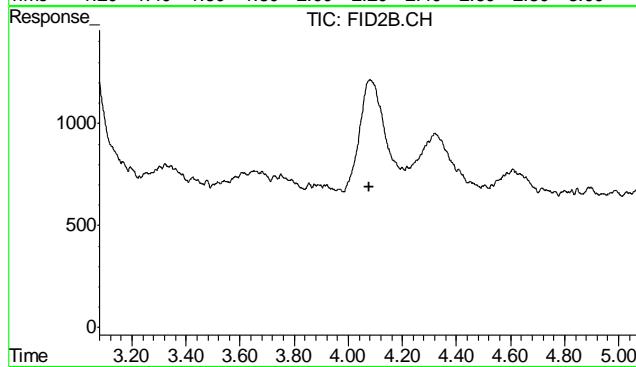
#2 1,2,4-Trichlorobenzene

R.T.: 14.336 min  
 Delta R.T.: -0.006 min  
 Response: 3279322  
 Conc: 108.71 % m



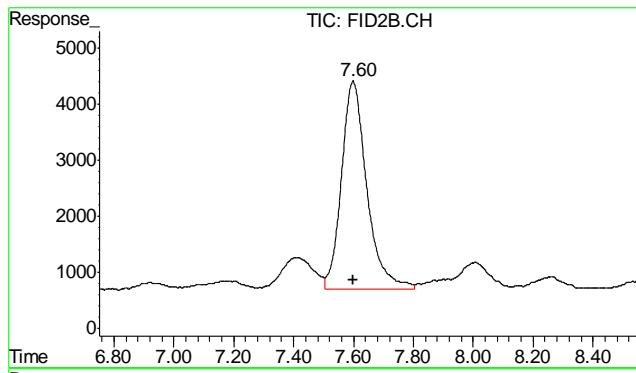
#4 Methyl-t-butyl-ether

R.T.: 0.000 min  
 Exp R.T. : 2.199 min  
 Response: 0  
 Conc: N.D.

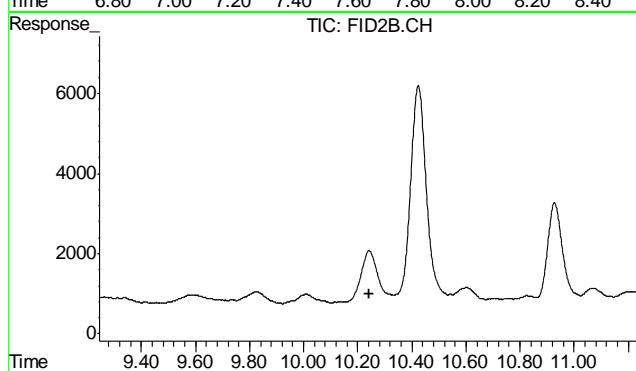


#5 Benzene

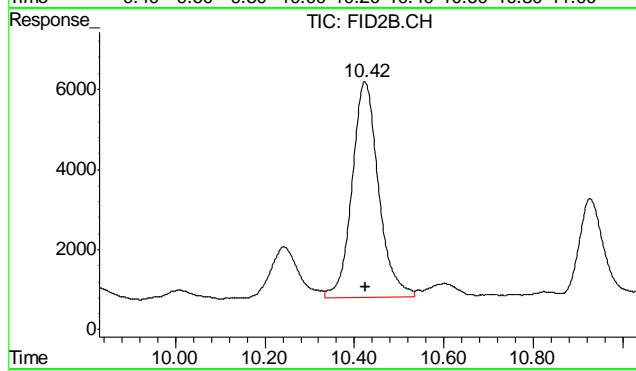
R.T.: 0.000 min  
 Exp R.T. : 4.079 min  
 Response: 0  
 Conc: N.D.



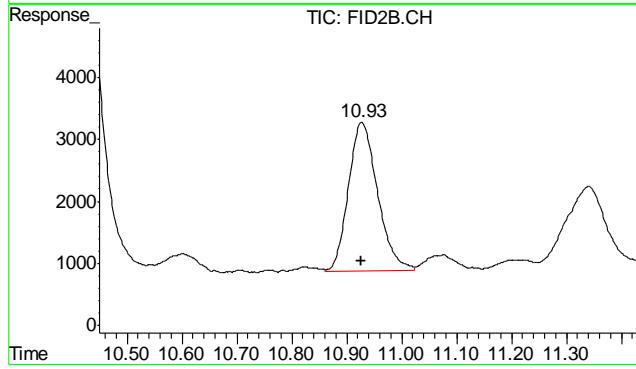
#6 Toluene  
R.T.: 7.599 min  
Delta R.T.: -0.002 min  
Response: 218482  
Conc: 0.40 ug/L



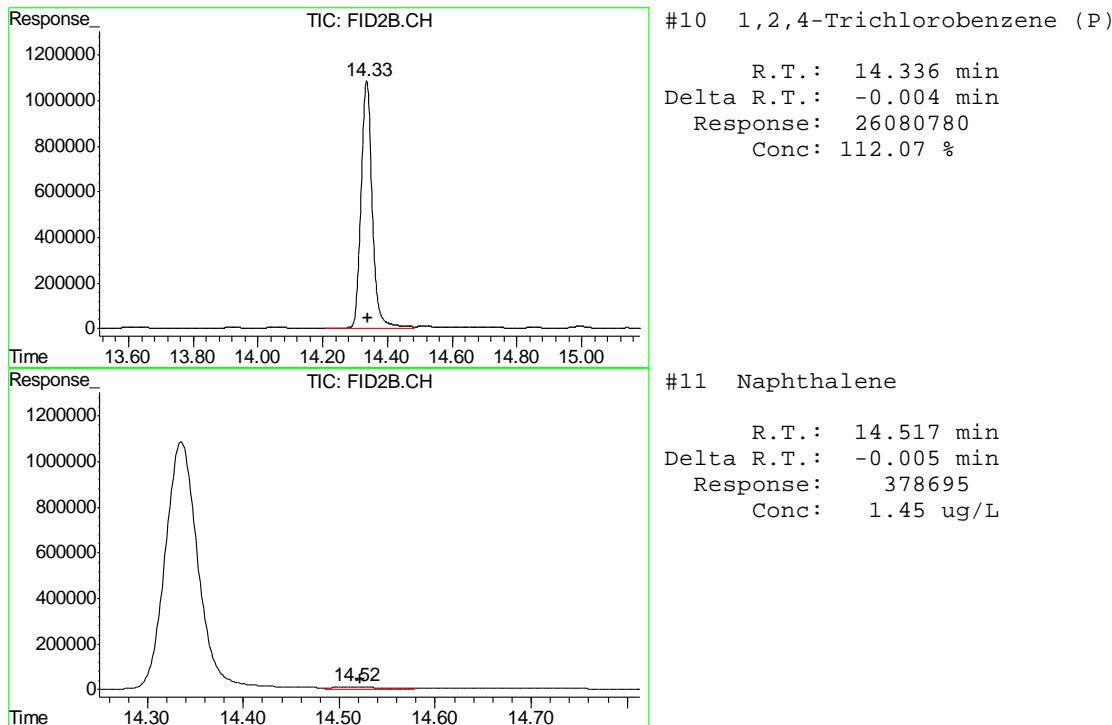
#7 Ethylbenzene  
R.T.: 0.000 min  
Exp R.T.: 10.244 min  
Response: 0  
Conc: N.D.



#8 m,p-Xylene  
R.T.: 10.424 min  
Delta R.T.: -0.003 min  
Response: 220790  
Conc: 0.39 ug/L



#9 o-Xylene  
R.T.: 10.928 min  
Delta R.T.: 0.000 min  
Response: 88601  
Conc: 0.19 ug/L

10.2.1  
**10**



## 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:** D32609  
**Account:** XTOKWR XTO Energy  
**Project:** FRU 297-32A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
OP5519-MB	FH002184.D 1		03/13/12	TR	03/12/12	OP5519	GFH113

The QC reported here applies to the following samples:

**Method:** SW846-8015B

D32609-1

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	88% 43-136%

11.11  
11

## Blank Spike Summary

Page 1 of 1

Job Number: D32609

Account: XTOKWR XTO Energy

Project: FRU 297-32A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
OP5519-BS	FH002186.D	1	03/13/12	TR	03/12/12	OP5519	GFH113

The QC reported here applies to the following samples:

Method: SW846-8015B

D32609-1

CAS No.	Compound	Spike mg/kg	BSP mg/kg	BSP %	Limits
	TPH-DRO (C10-C28)	667	536	80	58-130

CAS No.	Surrogate Recoveries	BSP	Limits
84-15-1	o-Terphenyl	86%	43-136%

11.2.1  
11

# Matrix Spike/Matrix Spike Duplicate Summary

Page 1 of 1

Job Number: D32609

Account: XTOKWR XTO Energy

Project: FRU 297-32A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
OP5519-MS	FH002216.D	50	03/14/12	TR	03/12/12	OP5519	GFH115
OP5519-MSD	FH002218.D	50	03/14/12	TR	03/12/12	OP5519	GFH115
D32595-1	FH002222.D	50	03/14/12	TR	03/12/12	OP5519	GFH115

The QC reported here applies to the following samples:

Method: SW846-8015B

D32609-1

CAS No.	Compound	D32595-1		Spike	MS	MS	MSD	MSD	RPD	Limits Rec/RPD
		mg/kg	Q	mg/kg	mg/kg	%	mg/kg	%		
	TPH-DRO (C10-C28)	43300		869	50300	805* a	41000	-264* a	20	20-183/43

CAS No.	Surrogate Recoveries	MS	MSD	D32595-1	Limits
84-15-1	o-Terphenyl	0% * b	0% * b	0% * b	43-136%

(a) Outside control limits due to high level in sample relative to spike amount.

(b) Outside control limits due to dilution.

11.3.1  
11



GC Semi-volatiles

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Raw Data

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## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\FH031312.SEC\  
 Data File : FH002202.D  
 Signal(s) : FID2B.ch  
 Acq On : 13 Mar 2012 9:20 pm  
 Operator : teder  
 Sample : D32609-1  
 Misc : OP5519,GFH113,30.01,,,2,1  
 ALS Vial : 62 Sample Multiplier: 1

Integration File: events.e  
 Quant Time: Mar 14 08:24:53 2012  
 Quant Method : C:\msdchem\1\METHODS\DRD-GFH95R.M  
 Quant Title : DRO-ORO REAR  
 QLast Update : Sun Mar 04 19:15:40 2012  
 Response via : Initial Calibration  
 Integrator: ChemStation

Volume Inj. :  
 Signal Phase :  
 Signal Info :

Compound	R.T.	Response	Conc	Units
<hr/>				
System Monitoring Compounds				
2) s o-Terphenyl	12.350	1106478169	646.171	ug/ml
<hr/>				
Target Compounds				
1) H TPH-DRO (C10-C28)	9.832	2519189305	1632.408	ug/ml
<hr/>				

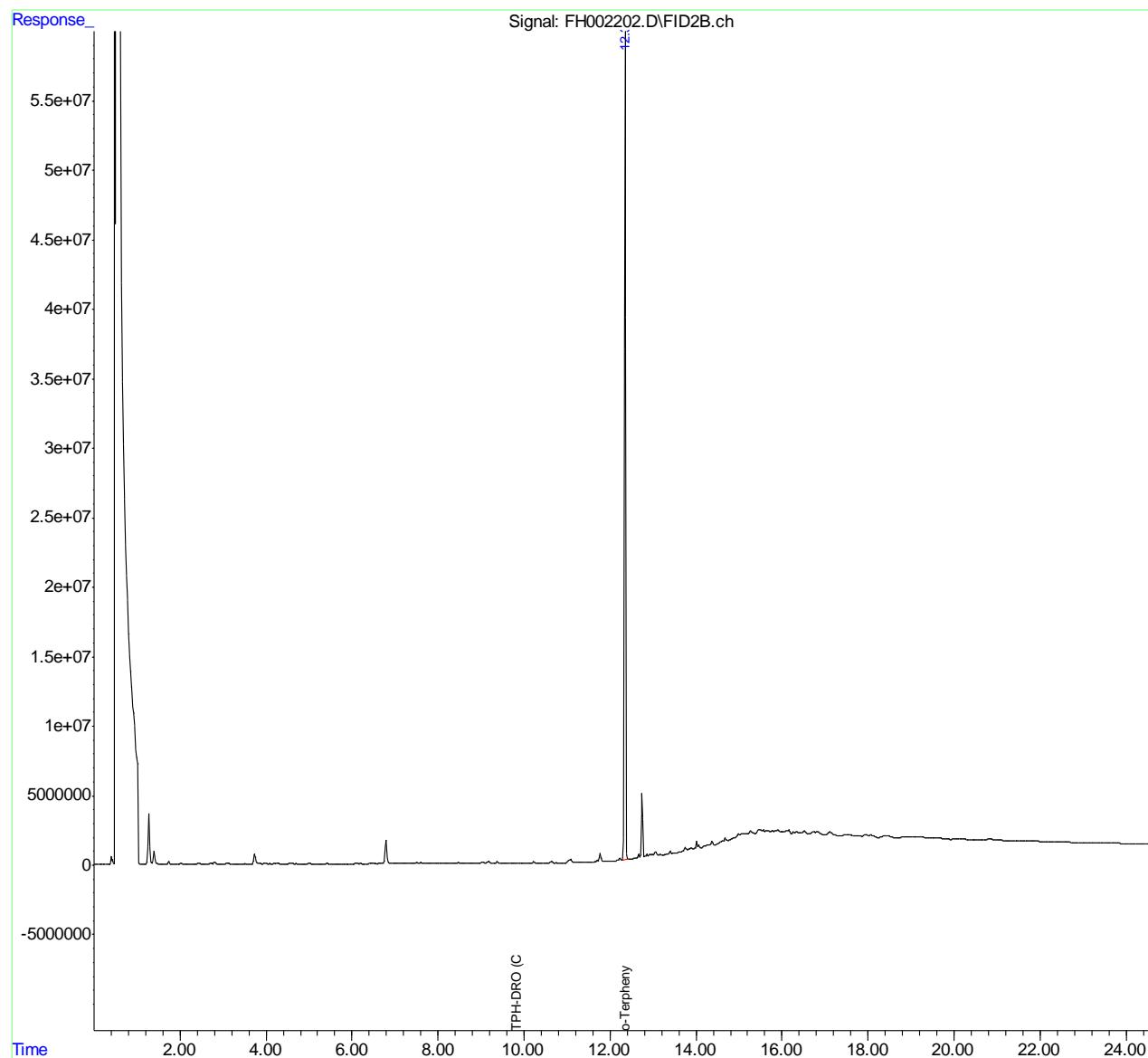
(f)=RT Delta > 1/2 Window (m)=manual int.

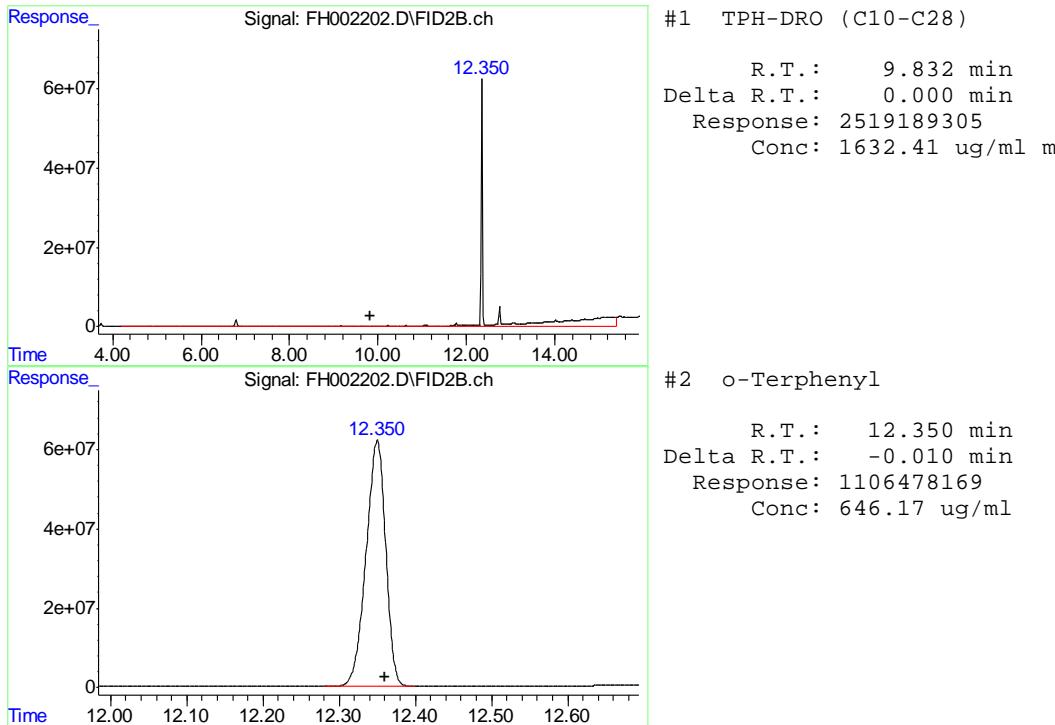
## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\FH031312.SEC\  
 Data File : FH002202.D  
 Signal(s) : FID2B.ch  
 Acq On : 13 Mar 2012 9:20 pm  
 Operator : tedr  
 Sample : D32609-1  
 Misc : OP5519,GFH113,30.01,,,2,1  
 ALS Vial : 62 Sample Multiplier: 1

Integration File: events.e  
 Quant Time: Mar 14 08:24:53 2012  
 Quant Method : C:\msdchem\1\METHODS\DRO-GFH95R.M  
 Quant Title : DRO-ORO REAR  
 QLast Update : Sun Mar 04 19:15:40 2012  
 Response via : Initial Calibration  
 Integrator: ChemStation

Volume Inj. :  
 Signal Phase :  
 Signal Info :





12.1.1

12

## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\FH031312.SEC\  
 Data File : FH002184.D  
 Signal(s) : FID2B.ch  
 Acq On : 13 Mar 2012 3:56 pm  
 Operator : teder  
 Sample : OP5519-MB  
 Misc : OP5519,GFH113,30.00,,,2,1  
 ALS Vial : 53 Sample Multiplier: 1

Integration File: events.e  
 Quant Time: Mar 13 17:12:48 2012  
 Quant Method : C:\msdchem\1\METHODS\YRO-GFH95R.M  
 Quant Title : DRO-ORO REAR  
 QLast Update : Sun Mar 04 19:15:40 2012  
 Response via : Initial Calibration  
 Integrator: ChemStation

Volume Inj. :  
 Signal Phase :  
 Signal Info :

Compound	R.T.	Response	Conc	Units
<hr/>				
System Monitoring Compounds				
2) s o-Terphenyl	12.356	1505589486	879.248	ug/ml
<hr/>				
Target Compounds				
1) H TPH-DRO (C10-C28)	9.832	75794120	49.114	ug/ml
<hr/>				

(f)=RT Delta &gt; 1/2 Window

(m)=manual int.

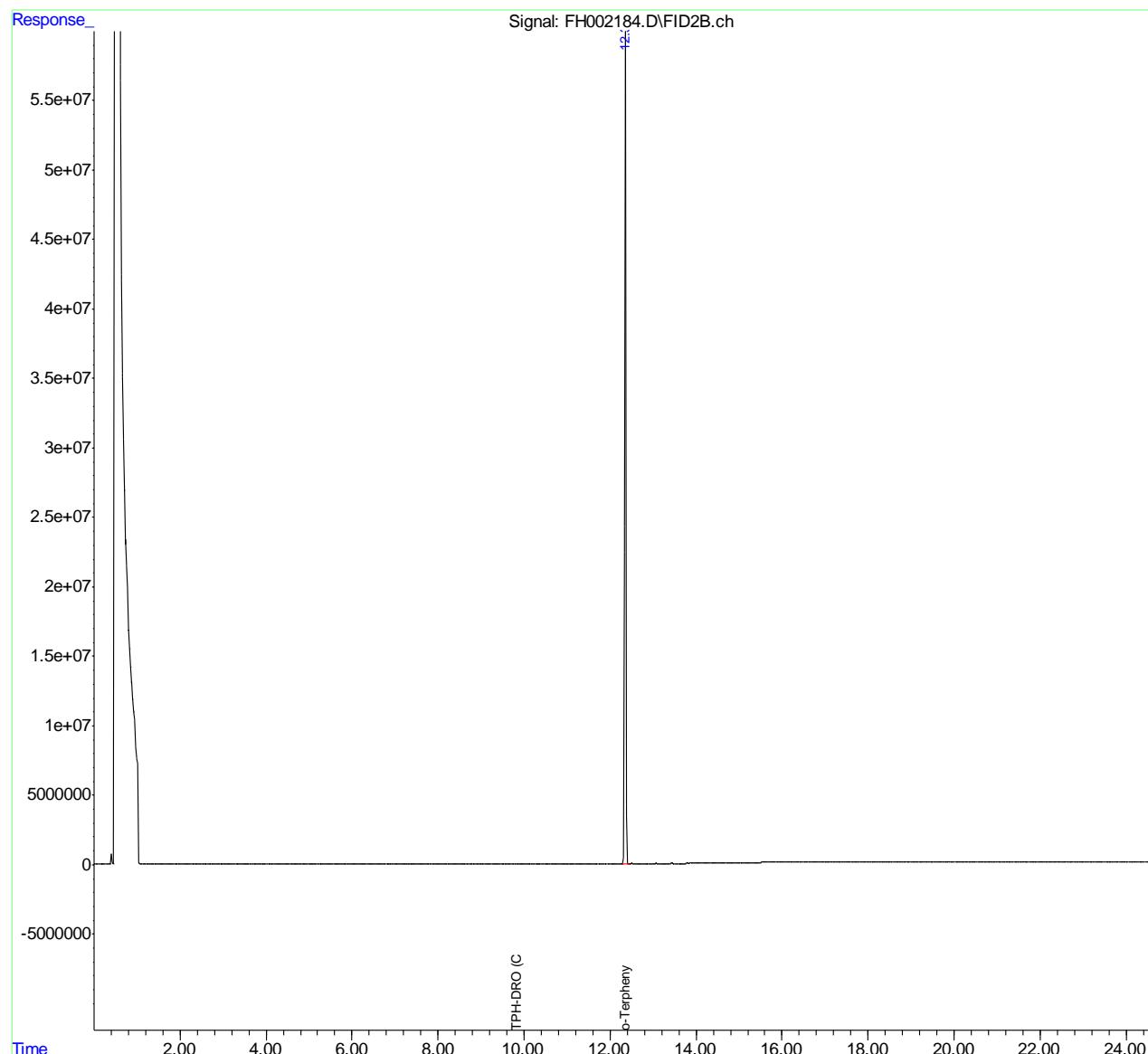
12.2.1  
12

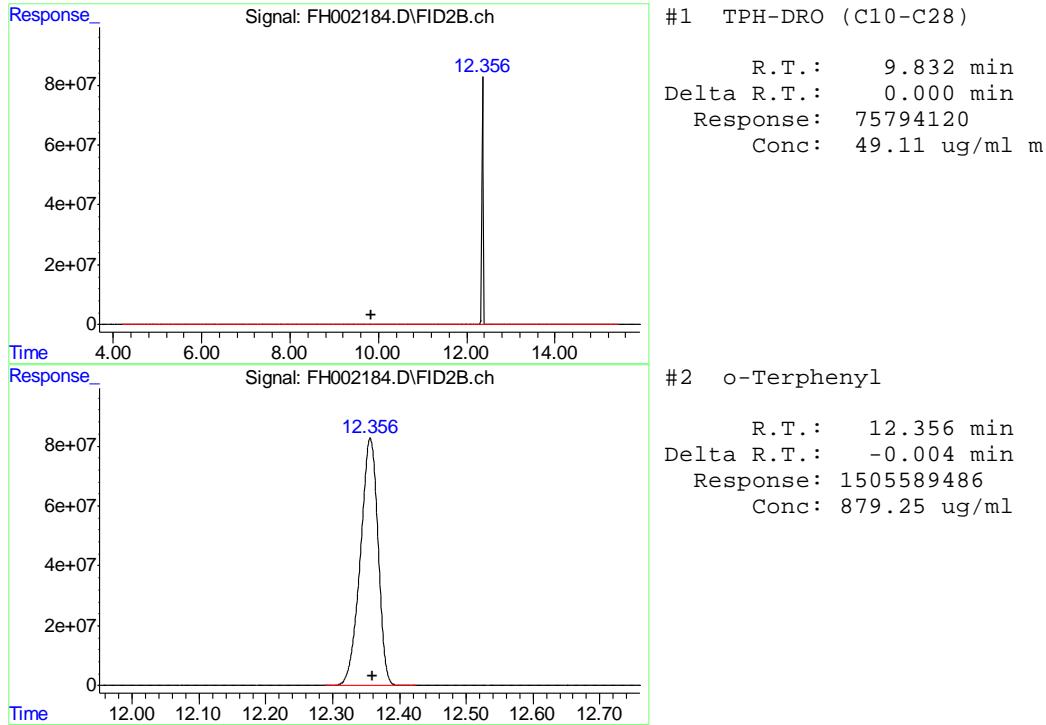
## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\FH031312.SEC\  
 Data File : FH002184.D  
 Signal(s) : FID2B.ch  
 Acq On : 13 Mar 2012 3:56 pm  
 Operator : tedr  
 Sample : OP5519-MB  
 Misc : OP5519,GFH113,30.00,,,2,1  
 ALS Vial : 53 Sample Multiplier: 1

Integration File: events.e  
 Quant Time: Mar 13 17:12:48 2012  
 Quant Method : C:\msdchem\1\METHODS\DRO-GFH95R.M  
 Quant Title : DRO-ORO REAR  
 QLast Update : Sun Mar 04 19:15:40 2012  
 Response via : Initial Calibration  
 Integrator: ChemStation

Volume Inj. :  
 Signal Phase :  
 Signal Info :





12.2.1

12



## 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: D32609  
Account: XTOKWR - XTO Energy  
Project: FRU 297-32A

QC Batch ID: MP7050  
Matrix Type: SOLID

Methods: SW846 7471B  
Units: mg/kg

Prep Date: 03/13/12

Metal	RL	IDL	MDL	MB raw	final
Mercury	0.10	.0011	.013	0.0020	<0.10

Associated samples MP7050: D32609-1

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: D32609  
Account: XTOKRWR - XTO Energy  
Project: FRU 297-32A

QC Batch ID: MP7050  
Matrix Type: SOLID

Methods: SW846 7471B  
Units: mg/kg

Prep Date: 03/13/12

Metal	D32609-1 Original MS	Spikelot HGWSR1	QC % Rec	QC Limits
Mercury	0.010	0.36	0.407	85.9 75-125

Associated samples MP7050: D32609-1

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: D32609  
Account: XTOKRWR - XTO Energy  
Project: FRU 297-32A

QC Batch ID: MP7050  
Matrix Type: SOLID

Methods: SW846 7471B  
Units: mg/kg

Prep Date:

03/13/12

Metal	D32609-1 Original	MSD	Spikelot HGWSR1	MSD % Rec	RPD	QC Limit
Mercury	0.010	0.37	0.407	88.4	2.7	

Associated samples MP7050: D32609-1

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: D32609  
Account: XTOKRWR - XTO Energy  
Project: FRU 297-32A

QC Batch ID: MP7050  
Matrix Type: SOLID

Methods: SW846 7471B  
Units: mg/kg

Prep Date: 03/13/12

Metal	BSP Result	Spikelot HGWSR1	QC % Rec	QC Limits
Mercury	0.40	0.4	100.0	80-120

Associated samples MP7050: D32609-1

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: D32609  
Account: XTOKRWR - XTO Energy  
Project: FRU 297-32A

QC Batch ID: MP7055  
Matrix Type: SOLID

Methods: SW846 6010C  
Units: mg/kg

Prep Date:

03/14/12

Metal	RL	IDL	MDL	MB raw	final
Aluminum	10	.59	.59		
Antimony	3.0	.31	.31		
Arsenic	2.5	.59	.59		
Barium	1.0	.11	.11	0.090	<1.0
Beryllium	1.0	.044	.1		
Boron	5.0	.48	.48		
Cadmium	1.0	.027	.27	0.010	<1.0
Calcium	40	.96	1.1		
Chromium	1.0	.018	.031	0.060	<1.0
Cobalt	0.50	.035	.035		
Copper	1.0	.085	.16	-0.18	<1.0
Iron	7.0	.34	2		
Lead	5.0	.16	.21	0.21	<5.0
Lithium	0.20	.028	.031		
Magnesium	20	.58	1.4		
Manganese	0.50	.0053	.012		
Molybdenum	1.0	.045	.054		
Nickel	3.0	.043	.099	0.0	<3.0
Phosphorus	10	1.1	1.2		
Potassium	200	5.5	9.2		
Selenium	5.0	.38	.5	-0.60	<5.0
Silicon	5.0	.38	.51		
Silver	3.0	.018	.051	-0.010	<3.0
Sodium	40	11	11		
Strontium	5.0		.017		
Thallium	1.0	.29	.34		
Tin	5.0	.55	1.3		
Titanium	1.0	.011	.1		
Uranium	5.0	.15	.2		
Vanadium	1.0	.016	.025		
Zinc	3.0	.028	.06	0.50	<3.0

Associated samples MP7055: D32609-1

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

BLANK RESULTS SUMMARY  
Part 2 - Method Blanks

Login Number: D32609  
Account: XTOKRWR - XTO Energy  
Project: FRU 297-32A

QC Batch ID: MP7055  
Matrix Type: SOLID

Methods: SW846 6010C  
Units: mg/kg

Prep Date:

Metal

(anr) Analyte not requested

## MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D32609  
 Account: XTOKRWR - XTO Energy  
 Project: FRU 297-32A

QC Batch ID: MP7055  
 Matrix Type: SOLID

Methods: SW846 6010C  
 Units: mg/kg

Prep Date: 03/14/12

Metal	D32609-1 Original MS	Spikelot MPICPALL	% Rec	QC Limits
Aluminum	anr			
Antimony	anr			
Arsenic	anr			
Barium	2100	2570	212	221.8(a) 75-125
Beryllium	anr			
Boron				
Cadmium	0.032	45.4	53	85.7 75-125
Calcium				
Chromium	42.6	82.3	53	75.0 75-125
Cobalt				
Copper	13.5	65.2	53	97.6 75-125
Iron	anr			
Lead	11.1	99.5	106	83.4 75-125
Lithium				
Magnesium	anr			
Manganese	anr			
Molybdenum	anr			
Nickel	19.8	61.7	53	79.1 75-125
Phosphorus	anr			
Potassium				
Selenium	0.52	90.0	106	84.5 75-125
Silicon				
Silver	0.13	19.7	21.2	92.4 75-125
Sodium				
Strontium				
Thallium	anr			
Tin				
Titanium				
Uranium	anr			
Vanadium				
Zinc	43.1	83.5	53	76.3 75-125

Associated samples MP7055: D32609-1

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

MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D32609  
Account: XTOKRWR - XTO Energy  
Project: FRU 297-32A

QC Batch ID: MP7055  
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.

## MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D32609  
 Account: XTOKRWR - XTO Energy  
 Project: FRU 297-32A

QC Batch ID: MP7055  
 Matrix Type: SOLID

Methods: SW846 6010C  
 Units: mg/kg

Prep Date:

03/14/12

Metal	D32609-1 Original	MSD	Spikelot MPICPALL	% Rec	MSD RPD	QC Limit
Aluminum	anr					
Antimony	anr					
Arsenic	anr					
Barium	2100	2690	216	272.9(a)	4.6	20
Beryllium	anr					
Boron						
Cadmium	0.032	46.6	54	86.2	2.6	20
Calcium						
Chromium	42.6	85.4	54	79.2	3.7	20
Cobalt						
Copper	13.5	66.5	54	98.1	2.0	20
Iron	anr					
Lead	11.1	99.9	108	82.2	0.4	20
Lithium						
Magnesium	anr					
Manganese	anr					
Molybdenum	anr					
Nickel	19.8	60.7	54	75.7	1.6	20
Phosphorus	anr					
Potassium						
Selenium	0.52	92.5	108	85.1	2.7	20
Silicon						
Silver	0.13	20.2	21.6	92.8	2.5	20
Sodium						
Strontium						
Thallium	anr					
Tin						
Titanium						
Uranium	anr					
Vanadium						
Zinc	43.1	110	54	101.2	4.2	20

Associated samples MP7055: D32609-1

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

MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D32609  
Account: XTOKRWR - XTO Energy  
Project: FRU 297-32A

QC Batch ID: MP7055  
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: D32609  
 Account: XTOKRWR - XTO Energy  
 Project: FRU 297-32A

QC Batch ID: MP7055  
 Matrix Type: SOLID

Methods: SW846 6010C  
 Units: mg/kg

Prep Date: 03/14/12

Metal	BSP Result	Spikelot MPICPALL	% Rec	QC Limits
Aluminum	anr			
Antimony	anr			
Arsenic	anr			
Barium	189	200	94.5	80-120
Beryllium	anr			
Boron				
Cadmium	47.2	50	94.4	80-120
Calcium				
Chromium	49.5	50	99.0	80-120
Cobalt				
Copper	47.5	50	95.0	80-120
Iron	anr			
Lead	96.9	100	96.9	80-120
Lithium				
Magnesium	anr			
Manganese	anr			
Molybdenum	anr			
Nickel	47.3	50	94.6	80-120
Phosphorus	anr			
Potassium				
Selenium	95.0	100	95.0	80-120
Silicon				
Silver	20.4	20	102.0	80-120
Sodium				
Strontium				
Thallium	anr			
Tin				
Titanium				
Uranium	anr			
Vanadium				
Zinc	47.6	50	95.2	80-120

Associated samples MP7055: D32609-1

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

SPIKE BLANK AND LAB CONTROL SAMPLE SUMMARY

Login Number: D32609  
Account: XTOKRWR - XTO Energy  
Project: FRU 297-32A

QC Batch ID: MP7055  
Matrix Type: SOLID

Methods: SW846 6010C  
Units: mg/kg

Prep Date:

Metal

(anr) Analyte not requested

13.2.3

13

## SERIAL DILUTION RESULTS SUMMARY

Login Number: D32609  
 Account: XTOKWR - XTO Energy  
 Project: FRU 297-32A

QC Batch ID: MP7055  
 Matrix Type: SOLID

Methods: SW846 6010C  
 Units: ug/l

Prep Date: 03/14/12

Metal	D32609-1 Original	SDL 1:5	%DIF	QC Limits
Aluminum	anr			
Antimony	anr			
Arsenic	anr			
Barium	22100	21300	9.0	0-10
Beryllium	anr			
Boron				
Cadmium	0.00	0.00	NC (a)	0-10
Calcium				
Chromium	458	462	16.0*(b)	0-10
Cobalt				
Copper	98.0	119	5.5	0-10
Iron	anr			
Lead	98.0	119	14.5*(b)	0-10
Lithium				
Magnesium	anr			
Manganese	anr			
Molybdenum	anr			
Nickel	234	224	20.8*(b)	0-10
Phosphorus	anr			
Potassium				
Selenium	0.00	0.00	NC (a)	0-10
Silicon				
Silver	4.00	3.00	150.0(a)	0-10
Sodium				
Strontium				
Thallium	anr			
Tin				
Titanium				
Uranium	anr			
Vanadium				
Zinc	517	506	25.4*(b)	0-10

Associated samples MP7055: D32609-1

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

SERIAL DILUTION RESULTS SUMMARY

Login Number: D32609  
Account: XTOKRWR - XTO Energy  
Project: FRU 297-32A

QC Batch ID: MP7055  
Matrix Type: SOLID

Methods: SW846 6010C  
Units: ug/l

Prep Date:

Metal

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

BLANK RESULTS SUMMARY  
Part 2 - Method Blanks

Login Number: D32609  
Account: XTOKRWR - XTO Energy  
Project: FRU 297-32A

QC Batch ID: MP7056  
Matrix Type: SOLID

Methods: SW846 6020A  
Units: mg/kg

Prep Date:

03/14/12

Metal	RL	IDL	MDL	MB raw	final
Aluminum	25	.14	1.3		
Antimony	0.20	.001	.012		
Arsenic	0.40	.049		0.036	<0.40
Barium	1.0	.0035	.025		
Beryllium	0.10	.0075	.055		
Boron	20	.97	.6		
Cadmium	0.050	.023	.034		
Calcium	200	1.8	9.5		
Chromium	1.0	.021	.041		
Cobalt	0.10	.0033	.0085		
Copper	1.0	.011	.055		
Iron	20	.81	18		
Lead	0.25	.0012	.023		
Magnesium	50	.067	.6		
Manganese	0.50	.007	.039		
Molybdenum	0.50	.0044	.025		
Nickel	1.0	.0029	.031		
Phosphorus	30	1.8	3.5		
Potassium	100	2	6		
Selenium	0.20	.075			
Silver	0.050	.0008	.022		
Sodium	250	.8	3		
Strontium	10	.004	.024		
Thallium	0.10	.015	.013		
Tin	5.0	.006	.15		
Titanium	1.0	.035	.12		
Uranium	0.25	.00038	.008		
Vanadium	2.0	.052	.19		
Zinc	5.0	.039	.23		

Associated samples MP7056: D32609-1

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

13.3.1  
13

## MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D32609  
 Account: XTOKRWR - XTO Energy  
 Project: FRU 297-32A

QC Batch ID: MP7056  
 Matrix Type: SOLID

Methods: SW846 6020A  
 Units: mg/kg

Prep Date: 03/14/12

Metal	D32609-1 Original MS	Spikelot MPICPALL	% Rec	QC Limits
Aluminum				
Antimony				
Arsenic	3.6	112	106	102.3 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 MP7056: D32609-1

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: D32609  
 Account: XTOKRWR - XTO Energy  
 Project: FRU 297-32A

QC Batch ID: MP7056  
 Matrix Type: SOLID

Methods: SW846 6020A  
 Units: mg/kg

Prep Date:

03/14/12

Metal	D32609-1 Original	MSD	Spikelot MPICPALL	% Rec	MSD RPD	QC Limit
Aluminum						
Antimony						
Arsenic	3.6	119	108	106.8	6.1	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 MP7056: D32609-1

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: D32609  
 Account: XTOKRWR - XTO Energy  
 Project: FRU 297-32A

QC Batch ID: MP7056  
 Matrix Type: SOLID

Methods: SW846 6020A  
 Units: mg/kg

Prep Date: 03/14/12

Metal	BSP Result	Spikelot MPICPALL	QC % Rec	Limits
Aluminum				
Antimony				
Arsenic	101	100	101.0	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 MP7056: D32609-1

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

13.3.3  
**13**

## SERIAL DILUTION RESULTS SUMMARY

Login Number: D32609  
 Account: XTOKWR - XTO Energy  
 Project: FRU 297-32A

QC Batch ID: MP7056  
 Matrix Type: SOLID

Methods: SW846 6020A  
 Units: ug/l

Prep Date: 03/14/12

Metal	D32609-1 Original	SDL 5:25	%DIF	QC Limits
Aluminum				
Antimony				
Arsenic	34.0	29.3	13.8*(a)	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 MP7056: D32609-1

Results < IDL are shown as zero for calculation purposes  
 (\*) Outside of QC limits  
 (anr) Analyte not requested  
 (a) Serial dilution indicates possible matrix interference.

BLANK RESULTS SUMMARY  
Part 2 - Method Blanks

Login Number: D32609  
Account: XTOKRWR - XTO Energy  
Project: FRU 297-32A

QC Batch ID: MP7061  
Matrix Type: AQUEOUS

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

Prep Date:

03/14/12

Metal	RL	IDL	MDL	MB raw	final
Aluminum	500	30	30		
Antimony	150	16	16		
Arsenic	130	30	30		
Barium	50	5.5	5.5		
Beryllium	50	2.2	2.5		
Boron	250	24	24		
Cadmium	50	1.4	1.4		
Calcium	2000	48	75	34.5	<2000
Chromium	50	.9	4		
Cobalt	25	1.8	1.8		
Copper	50	4.3	14		
Iron	350	17	65		
Lead	250	8	11		
Lithium	10	1.4	6		
Magnesium	1000	29	50	3.0	<1000
Manganese	25	.27	1.6		
Molybdenum	50	2.3	4.4		
Nickel	150	2.2	5		
Phosphorus	500	55	100		
Potassium	5000	280	280		
Selenium	250	19	19		
Silicon	250	19	19		
Silver	150	.9	1.6		
Sodium	2000	570	570	-240	<2000
Strontium	25		1.3		
Thallium	50	15	15		
Tin	250	28	50		
Titanium	50	.55	1.6		
Uranium	250	7.5	18		
Vanadium	50	.8	1.1		
Zinc	150	1.4	9		

Associated samples MP7061: D32609-1A

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

BLANK RESULTS SUMMARY  
Part 2 - Method Blanks

Login Number: D32609  
Account: XTOKRWR - XTO Energy  
Project: FRU 297-32A

QC Batch ID: MP7061  
Matrix Type: AQUEOUS

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

Prep Date:

Metal

(anr) Analyte not requested

13.4.1

13

## MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D32609  
 Account: XTOKRWR - XTO Energy  
 Project: FRU 297-32A

QC Batch ID: MP7061  
 Matrix Type: AQUEOUS

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

Prep Date: 03/14/12

Metal	D32609-1A Original MS	Spikelot MPICPALL	% Rec	QC Limits
Aluminum				
Antimony				
Arsenic				
Barium				
Beryllium				
Boron				
Cadmium				
Calcium	18000	154000	125000	108.8
Chromium				
Cobalt				
Copper				
Iron				
Lead				
Lithium				
Magnesium	14900	149000	125000	107.3
Manganese				
Molybdenum				
Nickel				
Phosphorus				
Potassium				
Selenium				
Silicon				
Silver				
Sodium	248000	392000	125000	115.2
Strontium				
Thallium				
Tin				
Titanium				
Uranium				
Vanadium				
Zinc				

Associated samples MP7061: D32609-1A

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

MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D32609  
Account: XTOKWR - XTO Energy  
Project: FRU 297-32A

QC Batch ID: MP7061  
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: D32609  
 Account: XTOKRWR - XTO Energy  
 Project: FRU 297-32A

QC Batch ID: MP7061  
 Matrix Type: AQUEOUS

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

Prep Date: 03/14/12

Metal	D32609-1A Original MSD	Spikelot MPICPALL	MSD % Rec	MSD RPD	QC Limit
Aluminum					
Antimony					
Arsenic					
Barium					
Beryllium					
Boron					
Cadmium					
Calcium	18000	156000	125000	110.4	1.3
Chromium					
Cobalt					
Copper					
Iron					
Lead					
Lithium					
Magnesium	14900	150000	125000	108.1	0.7
Manganese					
Molybdenum					
Nickel					
Phosphorus					
Potassium					
Selenium					
Silicon					
Silver					
Sodium	248000	396000	125000	118.4	1.0
Strontium					
Thallium					
Tin					
Titanium					
Uranium					
Vanadium					
Zinc					

Associated samples MP7061: D32609-1A

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

MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D32609  
Account: XTOKWR - XTO Energy  
Project: FRU 297-32A

QC Batch ID: MP7061  
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: D32609  
 Account: XTOKRWR - XTO Energy  
 Project: FRU 297-32A

QC Batch ID: MP7061  
 Matrix Type: AQUEOUS

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

Prep Date: 03/14/12

Metal	BSP Result	Spikelot MPICPALL	% Rec	QC Limits
Aluminum				
Antimony				
Arsenic				
Barium				
Beryllium				
Boron				
Cadmium				
Calcium	136000	125000	108.8	80-120
Chromium				
Cobalt				
Copper				
Iron				
Lead				
Lithium				
Magnesium	131000	125000	104.8	80-120
Manganese				
Molybdenum				
Nickel				
Phosphorus				
Potassium				
Selenium				
Silicon				
Silver				
Sodium	134000	125000	107.2	80-120
Strontium				
Thallium				
Tin				
Titanium				
Uranium				
Vanadium				
Zinc				

Associated samples MP7061: D32609-1A

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

SPIKE BLANK AND LAB CONTROL SAMPLE SUMMARY

Login Number: D32609  
Account: XTOKRWR - XTO Energy  
Project: FRU 297-32A

QC Batch ID: MP7061  
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: D32609  
Account: XTOKWR - XTO Energy  
Project: FRU 297-32A

Analyte	Batch ID	RL	MB Result	Units	Spike Amount	BSP Result	BSP %Recov	QC Limits
Chromium, Hexavalent	GP6713/GN14085	1.0	0.0	mg/kg	185	197	106.0	80-120%
Specific Conductivity	GP6711/GN14080			umhos/cm	9967	9850	98.8	90-110%
pH	GN14060			su	8.00	7.99	99.9	99.3-100.7%

Associated Samples:

Batch GN14060: D32609-1

Batch GP6711: D32609-1

Batch GP6713: D32609-1

(\*) Outside of QC limits

DUPLICATE RESULTS SUMMARY  
GENERAL CHEMISTRY

Login Number: D32609  
Account: XTOKWR - XTO Energy  
Project: FRU 297-32A

Analyte	Batch ID	QC Sample	Units	Original Result	DUP Result	RPD	QC Limits
Chromium, Hexavalent Redox Potential Vs H2	GP6713/GN14085 GN14069	D32493-1 D32609-1	mg/kg mv	0.0 391	0.0 398	0.0 1.8	0-20% 0-20%

Associated Samples:  
Batch GN14069: D32609-1  
Batch GP6713: D32609-1  
(\*) Outside of QC limits

MATRIX SPIKE RESULTS SUMMARY  
GENERAL CHEMISTRY

Login Number: D32609  
Account: XTOKWR - XTO Energy  
Project: FRU 297-32A

Analyte	Batch ID	QC Sample	Units	Original Result	Spike Amount	MS Result	%Rec	QC Limits
Chromium, Hexavalent	GP6713/GN14085	D32493-1	mg/kg	0.0	40	30.3	76.0	75-125%

Associated Samples:

Batch GP6713: D32609-1

(\*) Outside of QC limits

(N) Matrix Spike Rec. outside of QC limits

MATRIX SPIKE DUPLICATE RESULTS SUMMARY  
GENERAL CHEMISTRY

Login Number: D32609  
Account: XTOKWR - XTO Energy  
Project: FRU 297-32A

Analyte	Batch ID	QC Sample	Units	Original Result	Spike Amount	MSD Result	RPD	QC Limit
Chromium, Hexavalent	GP6713/GN14085	D32493-1	mg/kg	0.0	40	22.1		31.3(a)

Associated Samples:

Batch GP6713: D32609-1

(\*) Outside of QC limits

(N) Matrix Spike Rec. outside of QC limits

(a) Spike recovery indicates possible matrix interference and/or sample nonhomogeneity. Post digestion spike was run with a recovery of 117%.