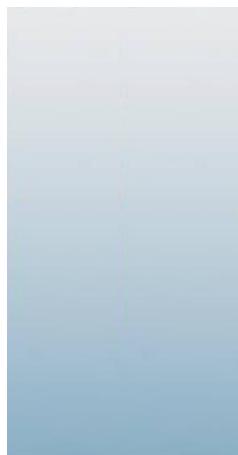




03/22/12



## Technical Report for

**XTO Energy**

**PCU T35X-2G**

**1108-11A**

**Accutest Job Number: D32747**

**Sampling Date: 03/14/12**

### Report to:

KRW Consulting, Inc.  
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ATTN: Dwayne Knudson

**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.

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

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

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

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: D32747

PCU T35X-2G

Project No: 1108-11A

Sample Number	Collected Date	Time By	Received	Matrix Code	Type	Client Sample ID
D32747-1	03/14/12	11:45 DS	03/15/12	SO	Soil	CUT SUBLINER
D32747-1A	03/14/12	11:45 DS	03/15/12	SO	Soil	CUT SUBLINER

---

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



## CASE NARRATIVE / CONFORMANCE SUMMARY

**Client:** XTO Energy

**Job No** D32747

**Site:** PCU T35X-2G

**Report Date** 3/22/2012 4:55:07 PM

On 03/15/2012, 1 sample(s), 0 Trip Blank(s), and 0 Field Blank(s) were received at Accutest Mountain States (AMS) at a temperature of 4 °C. The samples were intact and properly preserved, unless noted below. An AMS Job Number of D32747 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> V5V1216
------------------	--------------------------

- All samples were analyzed within the recommended method holding time.
- Sample(s) D32747-1MS, D32747-1MSD were used as the QC samples indicated.
- All method blanks for this batch meet method specific criteria.
- Sample(s) V5V1209-MB 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> OP5559
------------------	-------------------------

- All samples were extracted and analyzed within the recommended method holding time.
- All method blanks for this batch meet method specific criteria.
- Sample(s) D32747-1MS, D32747-1MSD were used as the QC samples indicated.

### Volatiles by GC By Method SW846 8015B

<b>Matrix</b> SO	<b>Batch ID:</b> GGB859
------------------	-------------------------

- All samples were analyzed within the recommended method holding time.
- All method blanks for this batch meet method specific criteria.
- Sample(s) D32747-1MS, D32747-1MSD were used as the QC samples indicated.

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

<b>Matrix</b> SO	<b>Batch ID:</b> OP5560
------------------	-------------------------

- All samples were extracted and analyzed within the recommended method holding time.
- All method blanks for this batch meet method specific criteria.
- Sample(s) D32747-1MS, D32747-1MSD were used as the QC samples indicated.

## Metals By Method SW846 6010C

**Matrix** AQ

**Batch ID:** MP7102

- All samples were digested and analyzed within the recommended method holding time.
- Sample(s) D32771-1AMS, D32771-1AMSD were used as the QC samples for the metals analysis.

**Matrix** SO

**Batch ID:** MP7078

- All samples were digested and analyzed within the recommended method holding time.
- All method blanks for this batch meet method specific criteria.
- Sample(s) D32747-1MS, D32747-1MSD, D32747-1SDL 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, Barium, Chromium, Nickel, Zinc are outside control limits for sample MP7078-SD1. Percent difference acceptable due to low initial sample concentration (< 50 times IDL).
- MP7078-SD1 for Barium, Chromium, Nickel, and Zinc: Serial dilution indicates possible matrix interference.

## Metals By Method SW846 6020A

**Matrix** SO

**Batch ID:** MP7079

- All samples were digested and analyzed within the recommended method holding time.
- All method blanks for this batch meet method specific criteria.
- Sample(s) D32747-1MS, D32747-1MSD, D32747-1SDL were used as the QC samples for the metals analysis.

## Metals By Method SW846 7471B

**Matrix** SO

**Batch ID:** MP7098

- All samples were digested and analyzed within the recommended method holding time.
- All method blanks for this batch meet method specific criteria.
- Sample(s) D32726-1MS, D32726-1MSD were used as the QC samples for the metals analysis.

## Wet Chemistry By Method ASTM D1498-76M

**Matrix** SO

**Batch ID:** GN14122

- Sample(s) D32747-1DUP were used as the QC samples for the Redox Potential Vs H<sub>2</sub> analysis.

## Wet Chemistry By Method DEPT.OF AG, BOOK N9

**Matrix** SO

**Batch ID:** GP6755

- All samples were prepared and analyzed within the recommended method holding time.
- All method blanks for this batch meet method specific criteria.

## Wet Chemistry By Method SM19 2540B M

**Matrix** SO

**Batch ID:** GN14097

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

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

**Matrix** SO

**Batch ID:** R12172

- The data for SW846 3060/7196A M meets quality control requirements.
- D32747-1 for Chromium, Trivalent: Calculated as: (Chromium) - (Chromium, Hexavalent)

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

**Matrix** SO

**Batch ID:** GP6737

- All samples were prepared and analyzed within the recommended method holding time.
- All method blanks for this batch meet method specific criteria.
- Sample(s) D32651-1DUP, D32651-1MS, D32651-1MSD were used as the QC samples for the Chromium, Hexavalent analysis.

## Wet Chemistry By Method SW846 9045C

**Matrix** SO

**Batch ID:** GN14113

- The following samples were run outside of holding time for method SW846 9045C: D32747-1

## Wet Chemistry By Method USDA HANDBOOK 60

**Matrix** SO

**Batch ID:** MP7102

- D32747-1A 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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Accutest Laboratories

**Report of Analysis**

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**Client Sample ID:** CUT SUBLINER  
**Lab Sample ID:** D32747-1  
**Matrix:** SO - Soil  
**Method:** SW846 8260B  
**Project:** PCU T35X-2G

**Date Sampled:** 03/14/12  
**Date Received:** 03/15/12  
**Percent Solids:** 93.8

	<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	5V20157.D	1	03/22/12	KV	n/a	n/a	V5V1216
Run #2							

	<b>Initial Weight</b>	<b>Final Volume</b>	<b>Methanol Aliquot</b>
Run #1	5.01 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	0.0761	0.057	0.025	mg/kg	
108-88-3	Toluene	0.226	0.11	0.057	mg/kg	
100-41-4	Ethylbenzene	0.0409	0.11	0.028	mg/kg	J
1330-20-7	Xylene (total)	0.296	0.23	0.11	mg/kg	

<b>CAS No.</b>	<b>Surrogate Recoveries</b>	<b>Run# 1</b>	<b>Run# 2</b>	<b>Limits</b>
2037-26-5	Toluene-D8	94%		61-130%
460-00-4	4-Bromofluorobenzene	88%		53-131%
17060-07-0	1,2-Dichloroethane-D4	89%		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 SUBLINER	<b>Date Sampled:</b>	03/14/12
<b>Lab Sample ID:</b>	D32747-1	<b>Date Received:</b>	03/15/12
<b>Matrix:</b>	SO - Soil	<b>Percent Solids:</b>	93.8
<b>Method:</b>	SW846 8270C BY SIM	SW846 3546	
<b>Project:</b>	PCU T35X-2G		
<b>File ID</b>	<b>DF</b>	<b>Analyzed</b>	<b>By</b>
Run #1	3G08570.D	1	03/19/12 DC
Run #2			
<b>Initial Weight</b>	<b>Final Volume</b>		
Run #1	30.0 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	0.0089	0.0046	mg/kg	
120-12-7	Anthracene	ND	0.0089	0.0046	mg/kg	
56-55-3	Benzo(a)anthracene	0.0168	0.0089	0.0046	mg/kg	
50-32-8	Benzo(a)pyrene	0.0084	0.0089	0.0046	mg/kg	J
205-99-2	Benzo(b)fluoranthene	0.0241	0.0089	0.0046	mg/kg	
207-08-9	Benzo(k)fluoranthene	ND	0.0089	0.0046	mg/kg	
218-01-9	Chrysene	0.0244	0.0089	0.0046	mg/kg	
53-70-3	Dibenz(a,h)anthracene	0.0058	0.0089	0.0046	mg/kg	J
206-44-0	Fluoranthene	0.0251	0.0089	0.0046	mg/kg	
86-73-7	Fluorene	ND	0.0089	0.0046	mg/kg	
193-39-5	Indeno(1,2,3-cd)pyrene	ND	0.0089	0.0046	mg/kg	
91-20-3	Naphthalene	0.0320	0.012	0.011	mg/kg	
129-00-0	Pyrene	0.0145	0.0089	0.0046	mg/kg	
CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits		
4165-60-0	Nitrobenzene-d5	50%		10-145%		
321-60-8	2-Fluorobiphenyl	44%		10-130%		
1718-51-0	Terphenyl-d14	70%		22-130%		

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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**Client Sample ID:** CUT SUBLINER  
**Lab Sample ID:** D32747-1  
**Matrix:** SO - Soil  
**Method:** SW846 8015B  
**Project:** PCU T35X-2G

**Date Sampled:** 03/14/12  
**Date Received:** 03/15/12  
**Percent Solids:** 93.8

	<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	GB15308.D	1	03/15/12	SK	n/a	n/a	GGB859
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)	ND	11	5.7	mg/kg	
<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	96%		60-140%		

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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<b>Client Sample ID:</b>	CUT SUBLINER	<b>Date Sampled:</b>	03/14/12
<b>Lab Sample ID:</b>	D32747-1	<b>Date Received:</b>	03/15/12
<b>Matrix:</b>	SO - Soil	<b>Percent Solids:</b>	93.8
<b>Method:</b>	SW846-8015B SW846 3546		
<b>Project:</b>	PCU T35X-2G		

	<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	FH002339.D	1	03/20/12	TR	03/19/12	OP5560	GFH123
Run #2							

	<b>Initial Weight</b>	<b>Final Volume</b>
Run #1	30.1 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)	31.1	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	64%		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 SUBLINER	<b>Date Sampled:</b>	03/14/12
<b>Lab Sample ID:</b>	D32747-1	<b>Date Received:</b>	03/15/12
<b>Matrix:</b>	SO - Soil	<b>Percent Solids:</b>	93.8
<b>Project:</b>	PCU T35X-2G		

**Metals Analysis**

Analyte	Result	RL	Units	DF	Prep	Analyzed By	Method	Prep Method
Arsenic	3.6	0.43	mg/kg	5	03/16/12	03/19/12 GJ	SW846 6020A <sup>2</sup>	SW846 3050B <sup>5</sup>
Barium	1880	1.1	mg/kg	1	03/16/12	03/16/12 JB	SW846 6010C <sup>1</sup>	SW846 3050B <sup>4</sup>
Cadmium	< 1.1	1.1	mg/kg	1	03/16/12	03/16/12 JB	SW846 6010C <sup>1</sup>	SW846 3050B <sup>4</sup>
Chromium	31.6	1.1	mg/kg	1	03/16/12	03/16/12 JB	SW846 6010C <sup>1</sup>	SW846 3050B <sup>4</sup>
Copper	15.5	1.1	mg/kg	1	03/16/12	03/16/12 JB	SW846 6010C <sup>1</sup>	SW846 3050B <sup>4</sup>
Lead	14.8	5.4	mg/kg	1	03/16/12	03/16/12 JB	SW846 6010C <sup>1</sup>	SW846 3050B <sup>4</sup>
Mercury	< 0.10	0.10	mg/kg	1	03/20/12	03/20/12 JB	SW846 7471B <sup>3</sup>	SW846 7471B <sup>6</sup>
Nickel	16.4	3.2	mg/kg	1	03/16/12	03/16/12 JB	SW846 6010C <sup>1</sup>	SW846 3050B <sup>4</sup>
Selenium	< 5.4	5.4	mg/kg	1	03/16/12	03/16/12 JB	SW846 6010C <sup>1</sup>	SW846 3050B <sup>4</sup>
Silver	< 3.2	3.2	mg/kg	1	03/16/12	03/16/12 JB	SW846 6010C <sup>1</sup>	SW846 3050B <sup>4</sup>
Zinc	48.2	3.2	mg/kg	1	03/16/12	03/16/12 JB	SW846 6010C <sup>1</sup>	SW846 3050B <sup>4</sup>

- (1) Instrument QC Batch: MA2266
- (2) Instrument QC Batch: MA2269
- (3) Instrument QC Batch: MA2272
- (4) Prep QC Batch: MP7078
- (5) Prep QC Batch: MP7079
- (6) Prep QC Batch: MP7098

RL = Reporting Limit

**Report of Analysis**

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**Client Sample ID:** CUT SUBLINER**Lab Sample ID:** D32747-1**Matrix:** SO - Soil**Date Sampled:** 03/14/12**Date Received:** 03/15/12**Percent Solids:** 93.8**Project:** PCU T35X-2G**General Chemistry**

Analyte	Result	RL	Units	DF	Analyzed	By	Method
Chromium, Hexavalent	< 1.0	1.0	mg/kg	1	03/20/12	CJ	SW846 3060A/7196A
Chromium, Trivalent <sup>a</sup>	31.6	2.1	mg/kg	1	03/20/12	CJ	SW846 3060/7196A M
Redox Potential Vs H2	377		mv	1	03/18/12 04:30	JK	ASTM D1498-76M
Solids, Percent	93.8		%	1	03/15/12	SWT	SM19 2540B M
Specific Conductivity	5480	1.0	umhos/cm	1	03/21/12	JK	DEPT.OF AG, BOOK N9
pH	9.35		su	1	03/16/12 10:45	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 SUBLINER	<b>Date Sampled:</b>	03/14/12
<b>Lab Sample ID:</b>	D32747-1A	<b>Date Received:</b>	03/15/12
<b>Matrix:</b>	SO - Soil	<b>Percent Solids:</b>	93.8
<b>Project:</b>	PCU T35X-2G		

**SAR Metals Analysis**

Analyte	Result	RL	Units	DF	Prep	Analyzed By	Method	Prep Method
Calcium	222	2.0	mg/l	1	03/20/12	03/20/12 JB	SW846 6010C <sup>1</sup>	EPA 200.7 <sup>2</sup>
Magnesium	64.4	1.0	mg/l	1	03/20/12	03/20/12 JB	SW846 6010C <sup>1</sup>	EPA 200.7 <sup>2</sup>
Sodium	1010	2.0	mg/l	1	03/20/12	03/20/12 JB	SW846 6010C <sup>1</sup>	EPA 200.7 <sup>2</sup>

(1) Instrument QC Batch: MA2270

(2) Prep QC Batch: MP7102

RL = Reporting Limit

**Report of Analysis**

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<b>Client Sample ID:</b>	CUT SUBLINER	<b>Date Sampled:</b>	03/14/12
<b>Lab Sample ID:</b>	D32747-1A	<b>Date Received:</b>	03/15/12
<b>Matrix:</b>	SO - Soil	<b>Percent Solids:</b>	93.8
<b>Project:</b>	PCU T35X-2G		

**General Chemistry**

Analyte	Result	RL	Units	DF	Analyzed	By	Method
Sodium Adsorption Ratio <sup>a</sup>	15.3		ratio	1	03/20/12 11:47	JB	USDA HANDBOOK 60

(a) Calculated as: (Na meq/L) / sqrt [(Ca meq/L)+ (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**

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

This  
one

D32747: Chain of Custody  
Page 1 of 2



## Accutest Laboratories Sample Receipt Summary

Accutest Job Number: D32747

Client: KRW

Immediate Client Services Action Required: No

Date / Time Received: 3/15/2012 12:00:00 PM

No. Coolers:

1

Client Service Action Required at Login: No

Project: XTO

Airbill #'s: 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:   | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| 2. Cooler temp verification: | Infrared gun                        |                          |
| 3. Cooler media:             | Ice (bag)                           |                          |

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

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

### Sample Integrity - Documentation

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

### Sample Integrity - Condition

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

### Sample Integrity - Instructions

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

Comments

Accutest Laboratories  
V:(303) 425-6021

4036 Youngfield Street  
F: (303) 425-6854

Wheat Ridge, CO  
www.accutest.com

4.1

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D32747: Chain of Custody

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## GC/MS Volatiles

5

### QC Data Summaries

Includes the following where applicable:

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

**Method Blank Summary**

**Job Number:** D32747  
**Account:** XTOKWR XTO Energy  
**Project:** PCU T35X-2G

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
V5V1216-MB	5V20148.D	1	03/22/12	KV	n/a	n/a	V5V1216

The QC reported here applies to the following samples:

**Method:** SW846 8260B

D32747-1

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	88%
17060-07-0	1,2-Dichloroethane-D4	96%

# Blank Spike Summary

Page 1 of 1

Job Number: D32747

Account: XTOKWR XTO Energy

Project: PCU T35X-2G

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
V5V1216-BS	5V20149.D	1	03/22/12	KV	n/a	n/a	V5V1216

The QC reported here applies to the following samples:

Method: SW846 8260B

D32747-1

CAS No.	Compound	Spike ug/kg	BSP ug/kg	BSP %	Limits
71-43-2	Benzene	50	50.9	102	70-130
100-41-4	Ethylbenzene	50	54.4	109	70-130
108-88-3	Toluene	50	52.2	104	70-130
1330-20-7	Xylene (total)	150	162	108	70-130

CAS No.	Surrogate Recoveries	BSP	Limits
2037-26-5	Toluene-D8	97%	61-130%
460-00-4	4-Bromofluorobenzene	92%	53-131%
17060-07-0	1,2-Dichloroethane-D4	81%	62-130%

# Matrix Spike/Matrix Spike Duplicate Summary

Page 1 of 1

Job Number: D32747

Account: XTOKWR XTO Energy

Project: PCU T35X-2G

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
D32747-1MS	5V20158.D	1	03/22/12	KV	n/a	n/a	V5V1216
D32747-1MSD	5V20159.D	1	03/22/12	KV	n/a	n/a	V5V1216
D32747-1	5V20157.D	1	03/22/12	KV	n/a	n/a	V5V1216

The QC reported here applies to the following samples:

Method: SW846 8260B

D32747-1

CAS No.	Compound	D32747-1		Spike	MS	MS	MSD	MSD	Limits	
		ug/kg	Q	ug/kg	ug/kg	%	ug/kg	%	RPD	Rec/RPD
71-43-2	Benzene	76.1		2830	2990	103	3080	106	3	70-134/30
100-41-4	Ethylbenzene	40.9	J	2830	3130	109	3200	112	2	70-137/30
108-88-3	Toluene	226		2830	3130	103	3150	103	1	70-130/30
1330-20-7	Xylene (total)	296		8480	9820	112	9880	113	1	61-131/30

CAS No.	Surrogate Recoveries	MS	MSD	D32747-1	Limits
2037-26-5	Toluene-D8	97%	96%	94%	61-130%
460-00-4	4-Bromofluorobenzene	104%	102%	88%	53-131%
17060-07-0	1,2-Dichloroethane-D4	85%	83%	89%	62-130%



## GC/MS Volatiles

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Raw Data

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## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\V5032212.S\  
 Data File : 5V20157.D  
 Acq On : 22 Mar 2012 2:35 pm  
 Operator : KOROUSHV  
 Sample : D32747-1  
 Misc : MS3611,V5V1216,5.007,,100,5,1  
 ALS Vial : 12 Sample Multiplier: 1

Quant Time: Mar 22 15:34:46 2012  
 Quant Method : C:\msdchem\1\METHODS\V5AP1212TVH1212.M  
 Quant Title : 8260  
 QLast Update : Wed Mar 21 09:50:04 2012  
 Response via : Initial Calibration

6.1.1

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
2) Pentafluorobenzene	11.647	168	263850	50.00	ug/l	0.00
35) 1,4-Difluorobenzene	12.446	114	403571	50.00	ug/l	0.00
53) Chlorobenzene-d5	15.095	117	447959	50.00	ug/l	0.00
74) 1,4-Dichlorobenzene-d4	17.070	152	281378	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	37529	44.51	ug/l	0.00
Spiked Amount 50.000	Range 70 - 130		Recovery	=	89.02%	
61) Toluene-d8	13.850	98	730358	46.92	ug/l	0.00
Spiked Amount 50.000	Range 70 - 130		Recovery	=	93.84%	
69) 4-Bromofluorobenzene	16.042	95	309347	44.12	ug/l	0.00
Spiked Amount 50.000	Range 70 - 130		Recovery	=	88.24%	

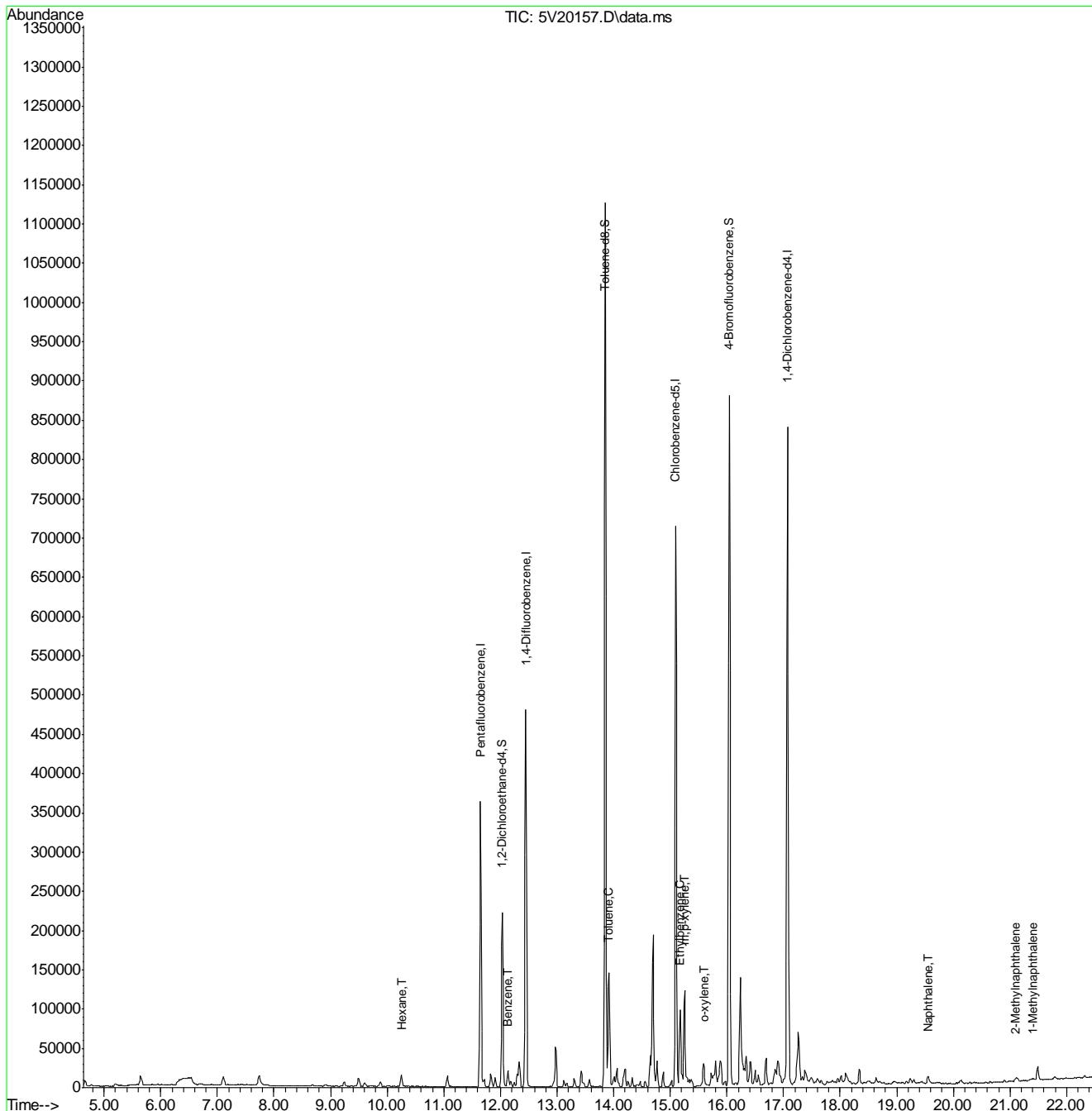
Target Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexane	10.254	57	7956	2.40	ug/l	100
50) Benzene	12.126	78	21559	1.35	ug/l	100
62) Toluene	13.908	92	46037	4.00	ug/l	100
66) Ethylbenzene	15.175	91	15384	0.72	ug/l	100
72) m,p-xylene	15.255	106	40909	4.63	ug/l	95
73) o-xylene	15.597	106	5051	0.61	ug/l	97
91) Naphthalene	19.559	128	7001	1.77	ug/l	100
94) 2-Methylnaphthalene	21.100	142	3864	1.32	ug/l	97
95) 1-Methylnaphthalene	21.397	142	1656	2.16	ug/l	94

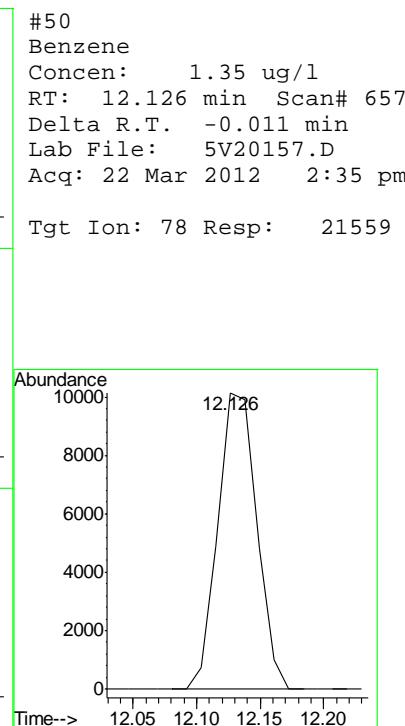
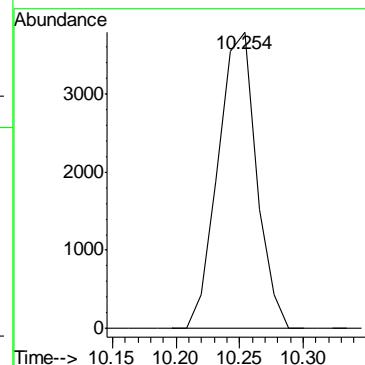
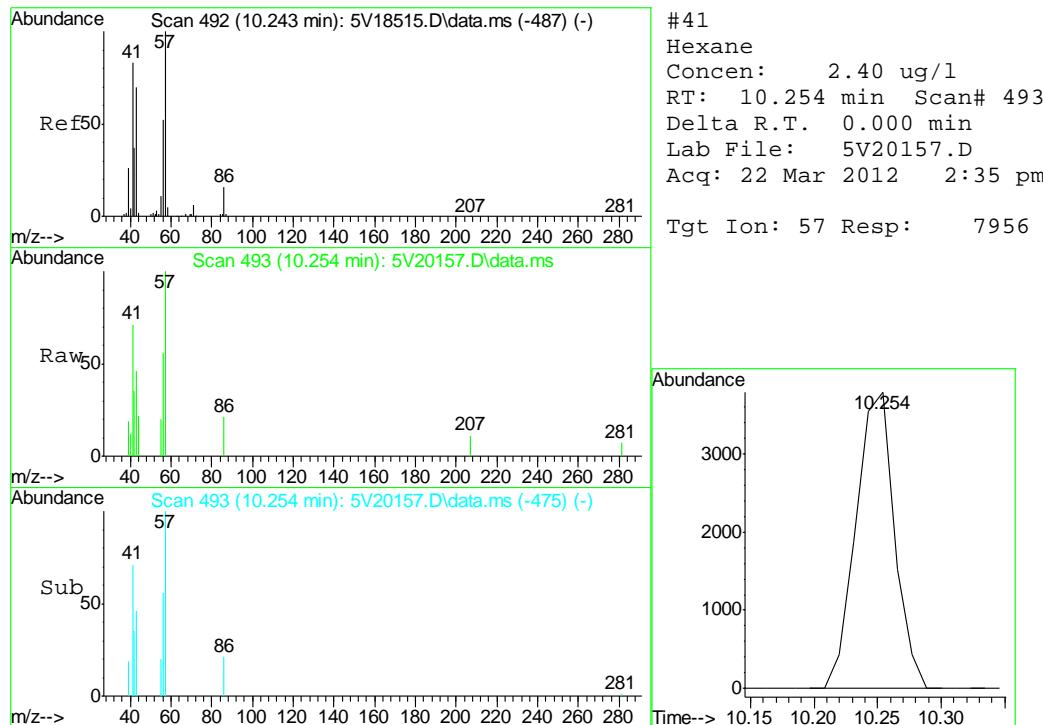
(#) = qualifier out of range (m) = manual integration (+) = signals summed

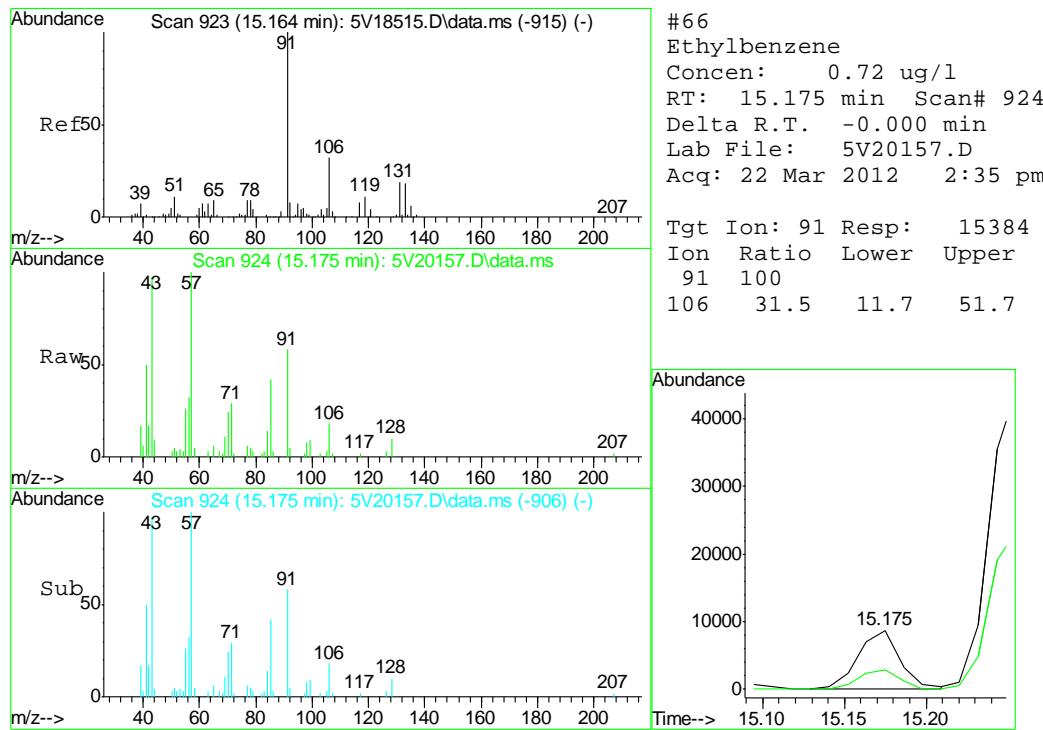
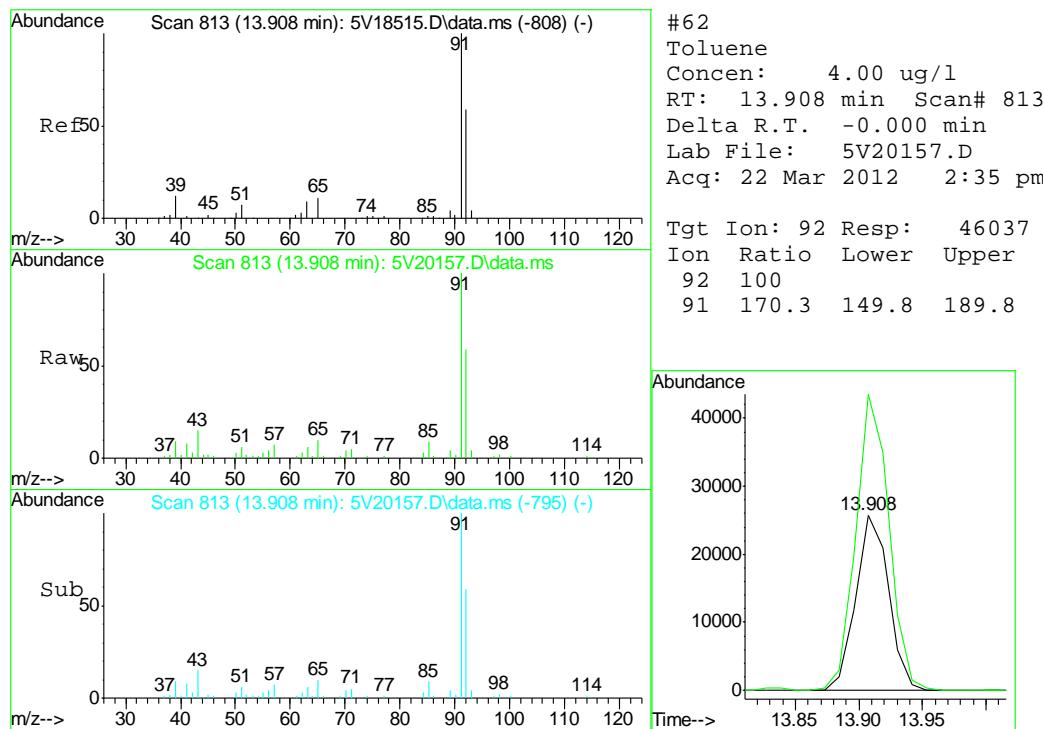
## Quantitation Report (QT Reviewed)

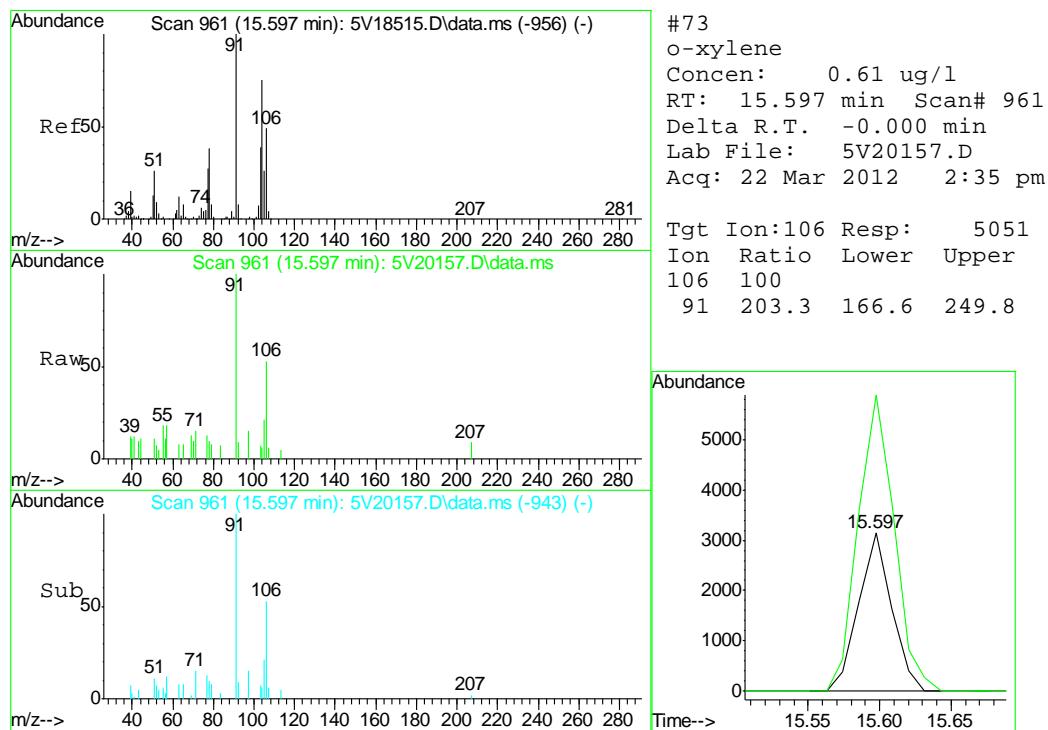
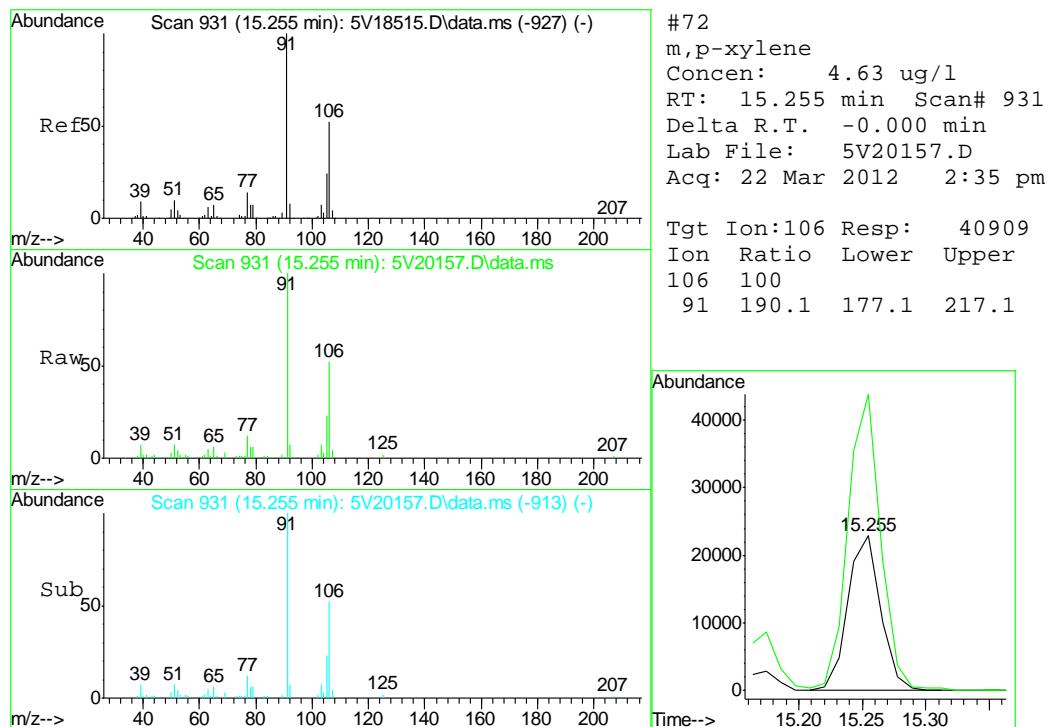
Data Path : C:\msdchem\1\DATA\V5032212.S\  
 Data File : 5V20157.D  
 Acq On : 22 Mar 2012 2:35 pm  
 Operator : KOROUSHV  
 Sample : D32747-1  
 Misc : MS3611,V5V1216,5.007,,100,5,1  
 ALS Vial : 12 Sample Multiplier: 1

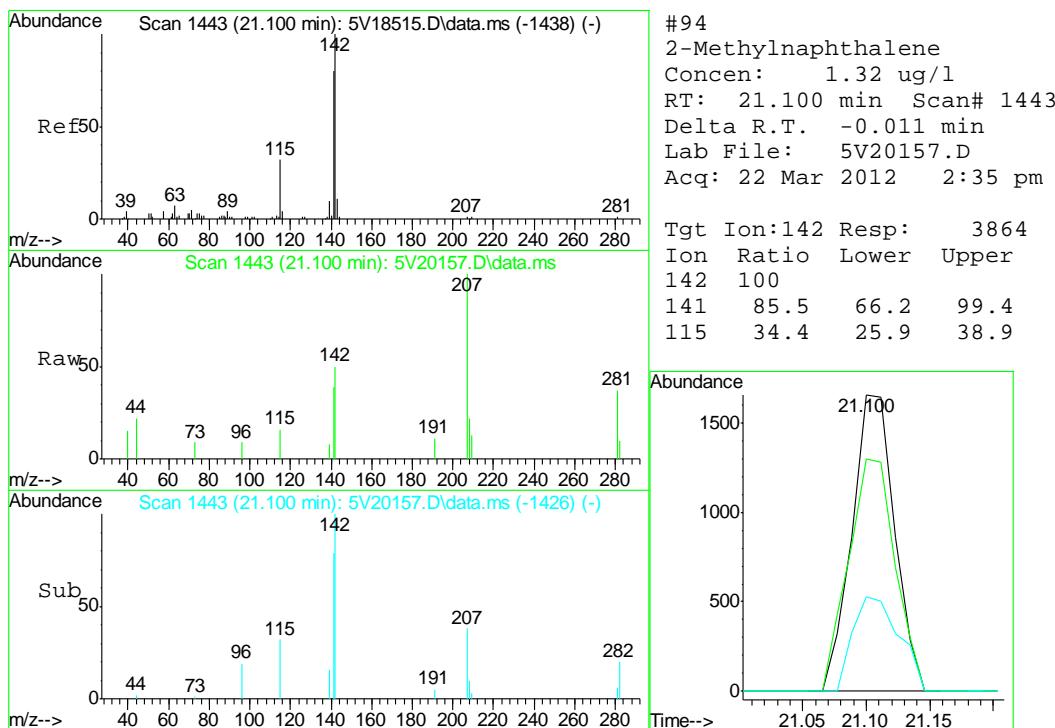
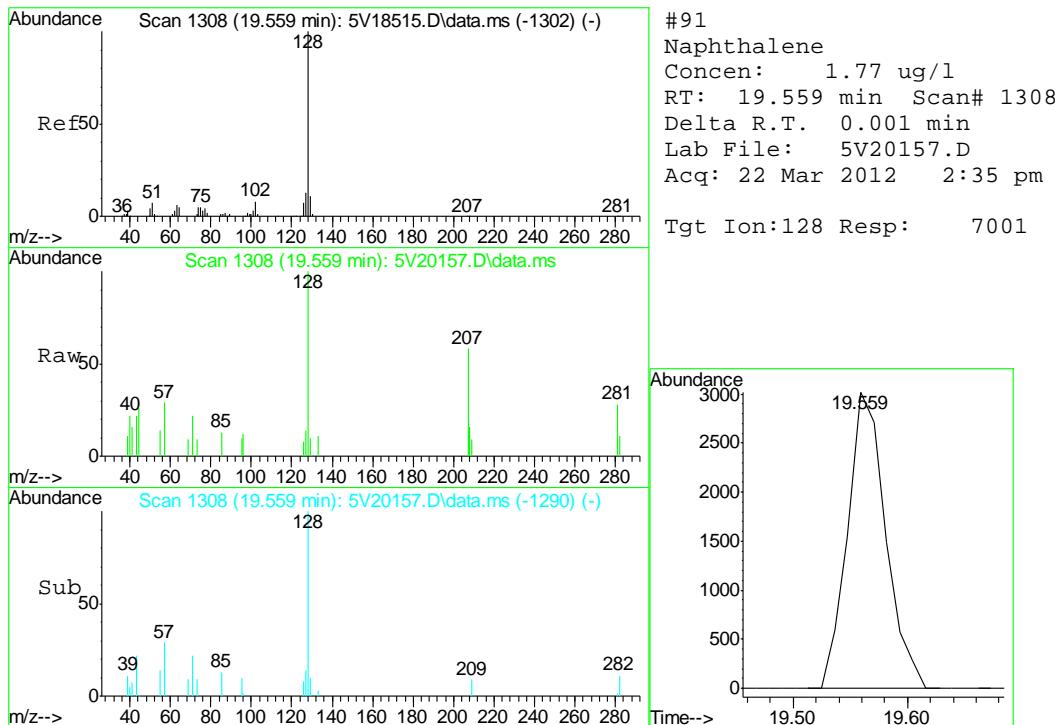
Quant Time: Mar 22 15:34:46 2012  
 Quant Method : C:\msdchem\1\METHODS\V5AP1212TVH1212.M  
 Quant Title : 8260  
 QLast Update : Wed Mar 21 09:50:04 2012  
 Response via : Initial Calibration

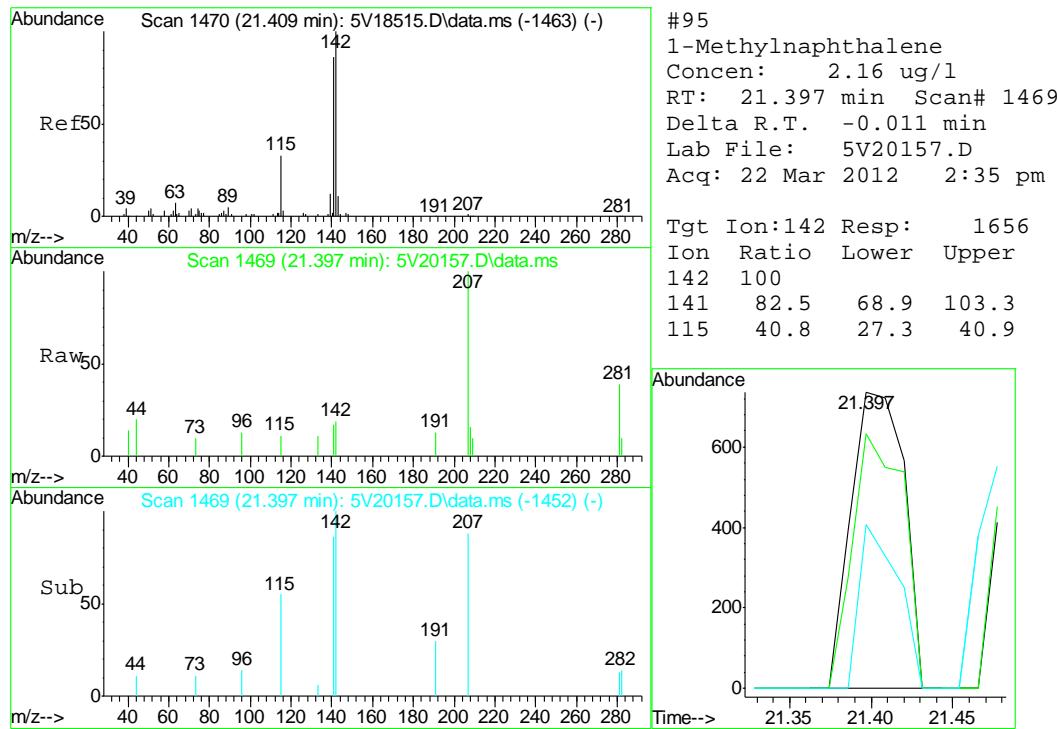












## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\V5032212.S\  
 Data File : 5V20148.D  
 Acq On : 22 Mar 2012 9:37 am  
 Operator : KOROUSHV  
 Sample : MB  
 Misc : MS3611,V5V1216,5.00,,100,5,1  
 ALS Vial : 3 Sample Multiplier: 1

Quant Time: Mar 22 13:49:12 2012  
 Quant Method : C:\msdchem\1\METHODS\V5AP1212TVH1212.M  
 Quant Title : 8260  
 QLast Update : Wed Mar 21 09:50:04 2012  
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
2) Pentafluorobenzene	11.647	168	300602	50.00	ug/l	0.00
35) 1,4-Difluorobenzene	12.446	114	455259	50.00	ug/l	0.00
53) Chlorobenzene-d5	15.095	117	471270	50.00	ug/l	0.00
74) 1,4-Dichlorobenzene-d4	17.070	152	278430	50.00	ug/l	0.00

System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	12.035	102	45796	48.02	ug/l	0.00
Spiked Amount	50.000	Range	70 - 130	Recovery	=	96.04%
61) Toluene-d8	13.850	98	850422	51.93	ug/l	0.00
Spiked Amount	50.000	Range	70 - 130	Recovery	=	103.86%
69) 4-Bromofluorobenzene	16.043	95	323804	43.90	ug/l	0.00
Spiked Amount	50.000	Range	70 - 130	Recovery	=	87.80%

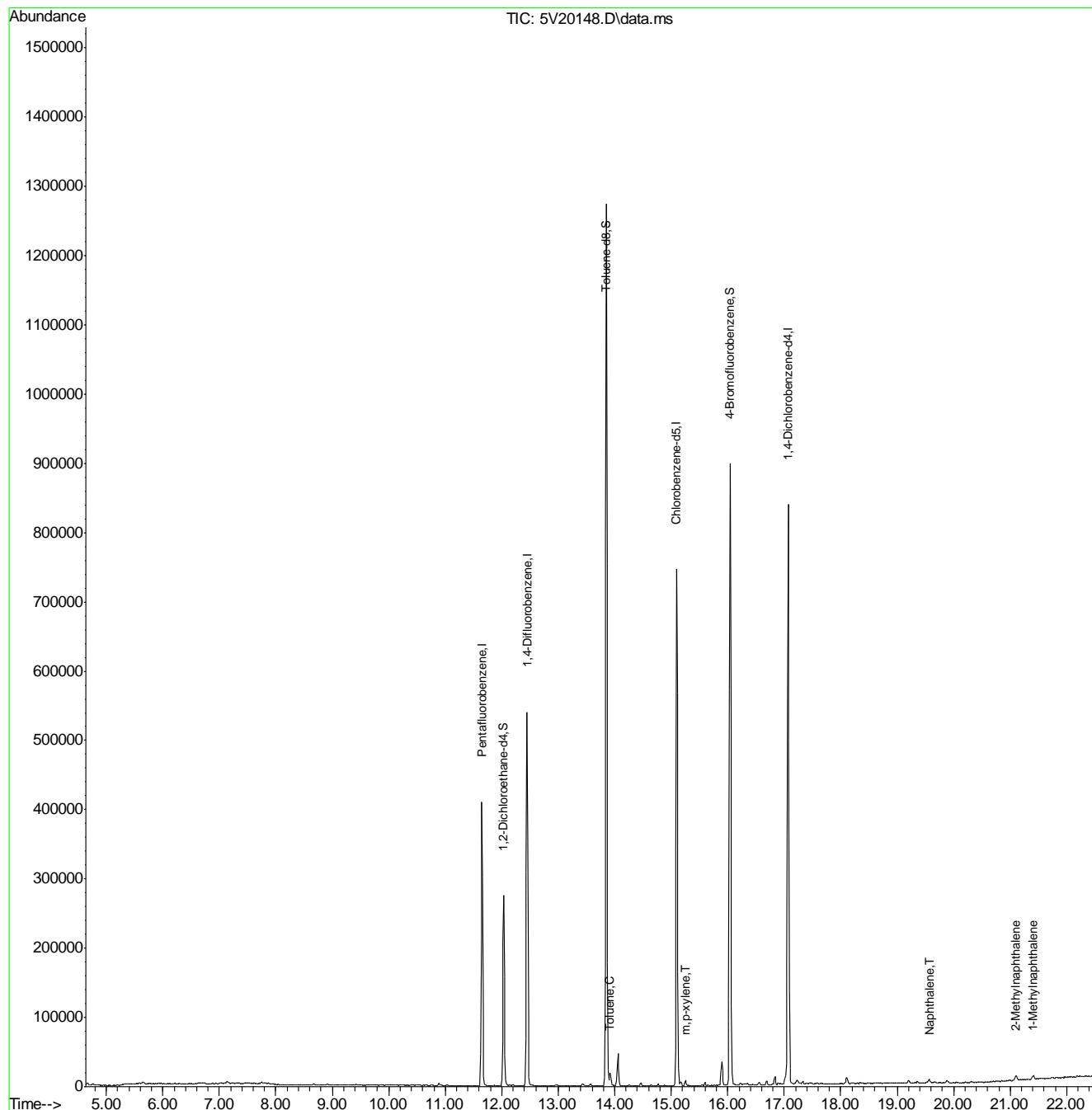
Target Compounds					Qvalue
62) Toluene	13.908	92	4209	0.35	ug/l
72) m,p-xylene	15.255	106	3324	0.36	ug/l
91) Naphthalene	19.570	128	8783	1.90	ug/l
94) 2-Methylnaphthalene	21.100	142	5886	2.02	ug/l
95) 1-Methylnaphthalene	21.408	142	5257	3.05	ug/l

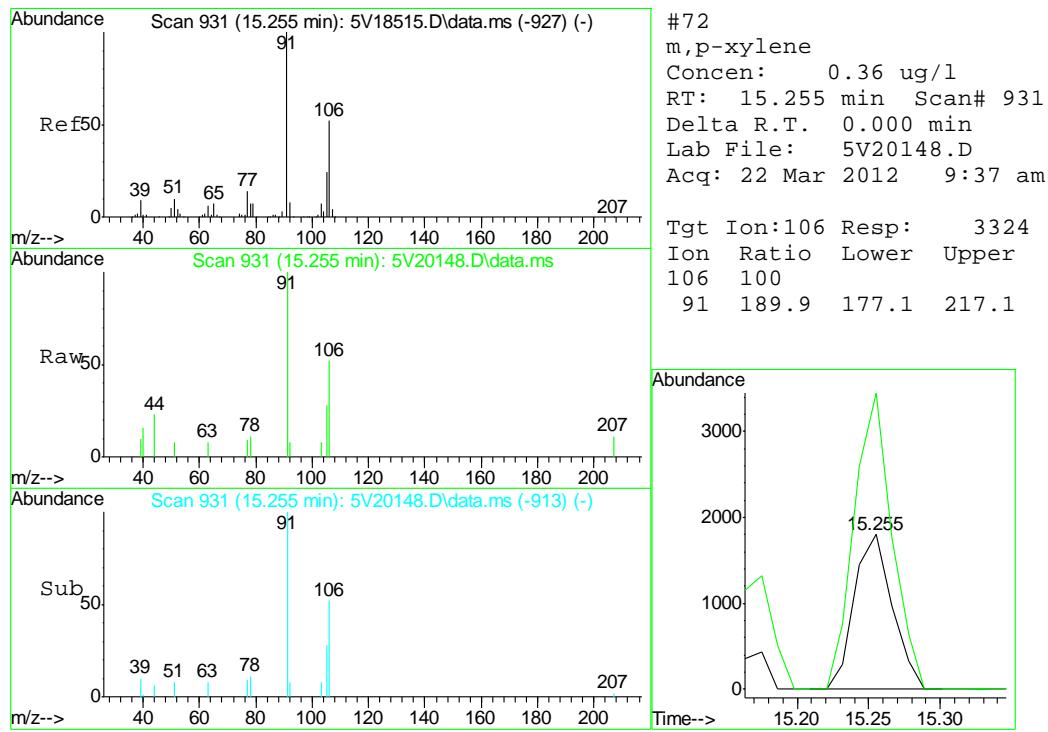
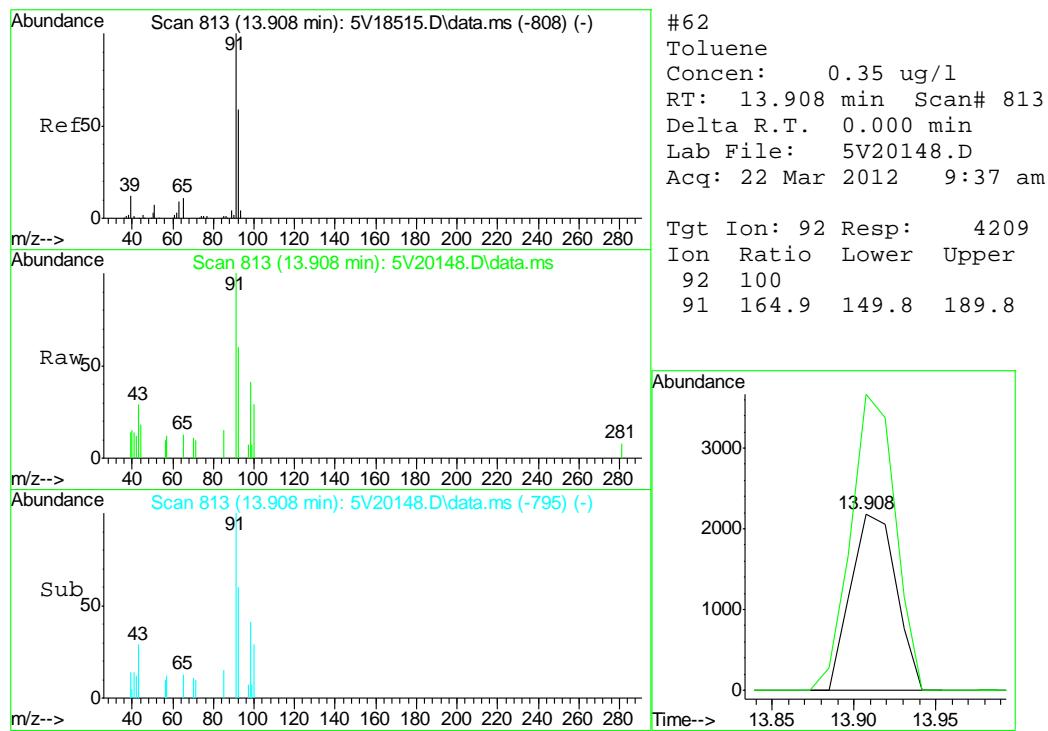
(#) = qualifier out of range (m) = manual integration (+) = signals summed

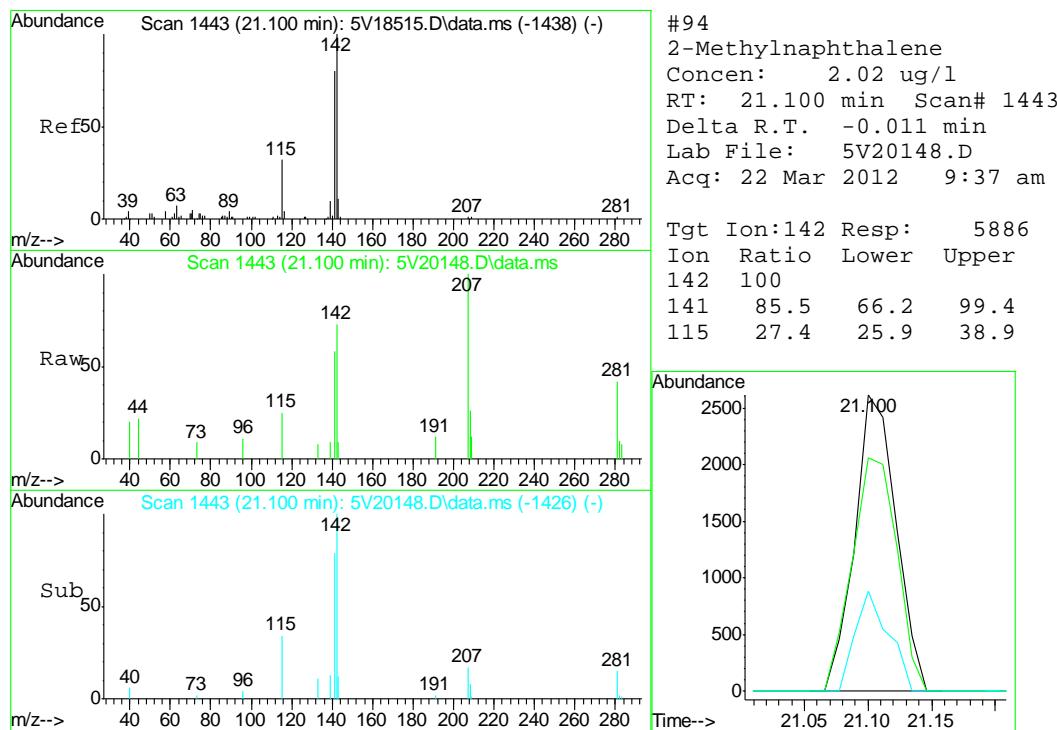
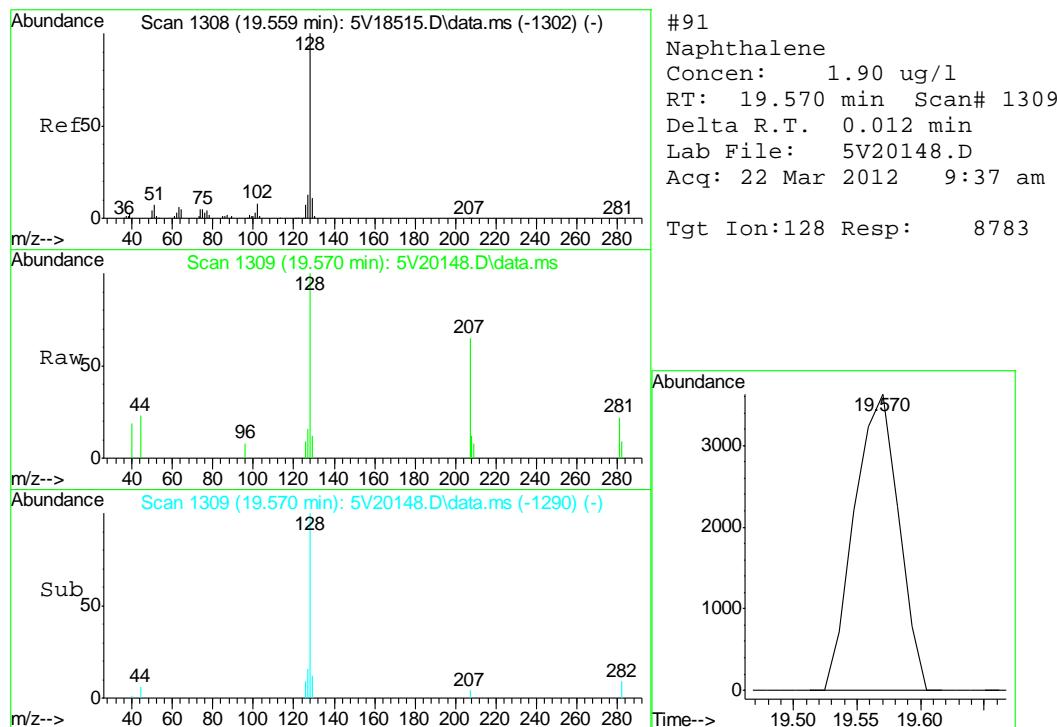
## Quantitation Report (QT Reviewed)

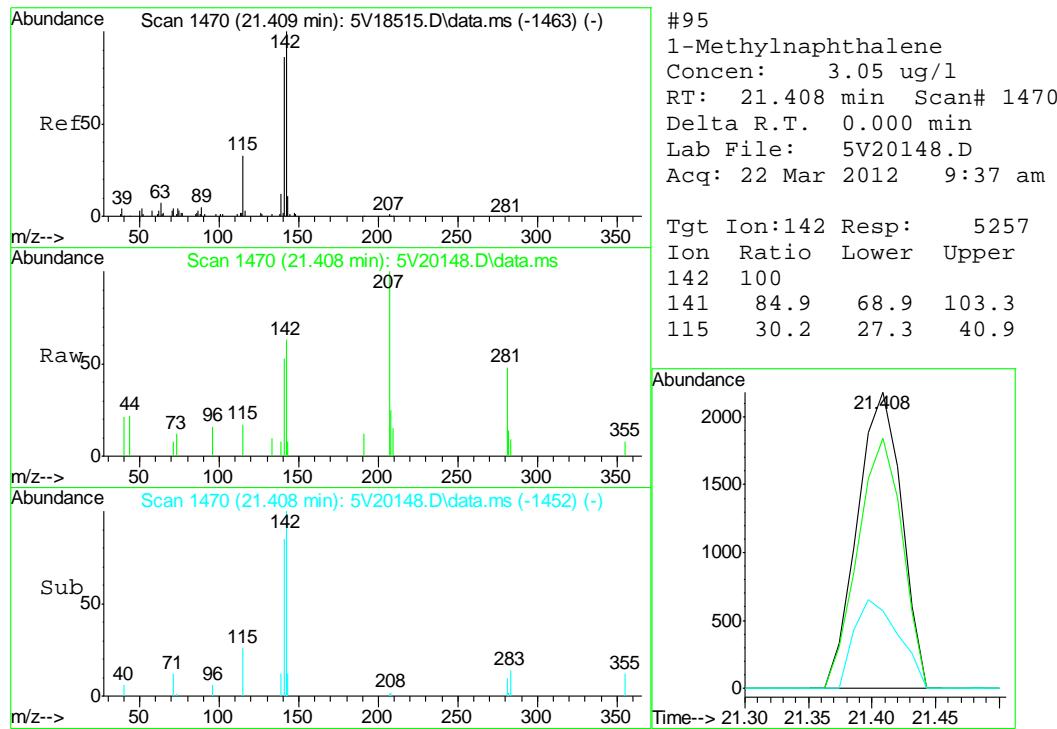
Data Path : C:\msdchem\1\DATA\V5032212.S\  
 Data File : 5V20148.D  
 Acq On : 22 Mar 2012 9:37 am  
 Operator : KOROUSHV  
 Sample : MB  
 Misc : MS3611,V5V1216,5.00,,100,5,1  
 ALS Vial : 3 Sample Multiplier: 1

Quant Time: Mar 22 13:49:12 2012  
 Quant Method : C:\msdchem\1\METHODS\V5AP1212TVH1212.M  
 Quant Title : 8260  
 QLast Update : Wed Mar 21 09:50:04 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: D32747

Account: XTOKWR XTO Energy

Project: PCU T35X-2G

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
OP5559-MB	3G08568.D	1	03/19/12	DC	03/19/12	OP5559	E3G352

The QC reported here applies to the following samples:

**Method:** SW846 8270C BY SIM

D32747-1

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	89%
321-60-8	2-Fluorobiphenyl	71%
1718-51-0	Terphenyl-d14	90%

## Blank Spike Summary

Page 1 of 1

Job Number: D32747

Account: XTOKWR XTO Energy

Project: PCU T35X-2G

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
OP5559-BS	3G08569.D	1	03/19/12	DC	03/19/12	OP5559	E3G352

The QC reported here applies to the following samples:

Method: SW846 8270C BY SIM

D32747-1

CAS No.	Compound	Spike ug/kg	BSP ug/kg	BSP %	Limits
83-32-9	Acenaphthene	83.3	72.6	87	34-130
120-12-7	Anthracene	83.3	76.6	92	35-130
56-55-3	Benzo(a)anthracene	83.3	69.5	83	36-130
50-32-8	Benzo(a)pyrene	83.3	67.7	81	36-130
205-99-2	Benzo(b)fluoranthene	83.3	66.3	80	35-130
207-08-9	Benzo(k)fluoranthene	83.3	71.3	86	37-130
218-01-9	Chrysene	83.3	73.0	88	40-130
53-70-3	Dibenzo(a,h)anthracene	83.3	69.6	84	32-130
206-44-0	Fluoranthene	83.3	77.6	93	38-130
86-73-7	Fluorene	83.3	75.8	91	35-130
193-39-5	Indeno(1,2,3-cd)pyrene	83.3	75.8	91	28-130
91-20-3	Naphthalene	83.3	72.4	87	35-130
129-00-0	Pyrene	83.3	74.0	89	37-130

CAS No.	Surrogate Recoveries	BSP	Limits
4165-60-0	Nitrobenzene-d5	91%	10-145%
321-60-8	2-Fluorobiphenyl	71%	10-130%
1718-51-0	Terphenyl-d14	83%	22-130%

# Matrix Spike/Matrix Spike Duplicate Summary

Page 1 of 1

Job Number: D32747

Account: XTOKWR XTO Energy

Project: PCU T35X-2G

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
OP5559-MS	3G08571.D	1	03/19/12	DC	03/19/12	OP5559	E3G352
OP5559-MSD	3G08572.D	1	03/19/12	DC	03/19/12	OP5559	E3G352
D32747-1	3G08570.D	1	03/19/12	DC	03/19/12	OP5559	E3G352

The QC reported here applies to the following samples:

Method: SW846 8270C BY SIM

D32747-1

7.3.1

CAS No.	Compound	D32747-1		Spike	MS	MS	MSD	MSD	RPD	Limits Rec/RPD
		ug/kg	Q	ug/kg	ug/kg	%	ug/kg	%		
83-32-9	Acenaphthene	ND		88.8	60.3	68	58.9	66	2	10-155/30
120-12-7	Anthracene	ND		88.8	76.3	86	71.2	80	7	10-155/30
56-55-3	Benzo(a)anthracene	16.8		88.8	94.3	87	89.8	82	5	10-175/30
50-32-8	Benzo(a)pyrene	8.4	J	88.8	76.1	76	71.4	71	6	10-164/30
205-99-2	Benzo(b)fluoranthene	24.1		88.8	104	90	102	88	2	10-165/30
207-08-9	Benzo(k)fluoranthene	ND		88.8	62.3	70	55.5	63	12	10-178/30
218-01-9	Chrysene	24.4		88.8	97.2	82	93.0	77	4	10-147/30
53-70-3	Dibenzo(a,h)anthracene	5.8	J	88.8	99.3	105	94.1	100	5	10-144/30
206-44-0	Fluoranthene	25.1		88.8	94.6	78	87.9	71	7	10-207/30
86-73-7	Fluorene	ND		88.8	77.4	87	75.1	85	3	10-163/30
193-39-5	Indeno(1,2,3-cd)pyrene	ND		88.8	126	142	121	137	4	10-180/30
91-20-3	Naphthalene	32.0		88.8	88.1	63	95.9	72	8	10-198/30
129-00-0	Pyrene	14.5		88.8	104	101	98.3	95	6	10-189/30

CAS No.	Surrogate Recoveries	MS	MSD	D32747-1	Limits
4165-60-0	Nitrobenzene-d5	58%	63%	50%	10-145%
321-60-8	2-Fluorobiphenyl	54%	57%	44%	10-130%
1718-51-0	Terphenyl-d14	81%	82%	70%	22-130%



## GC/MS Semi-volatiles

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Raw Data

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Judy Nelson  
 03/20/12 13:45

## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\031912\  
 Data File : 3g08570.D  
 Acq On : 19 Mar 2012 1:38 pm  
 Operator : DONC  
 Sample : D32747-1  
 Misc : OP5559,E3G352,30.02,,,1,1  
 ALS Vial : 6 Sample Multiplier: 1

Quant Time: Mar 20 10:02:52 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

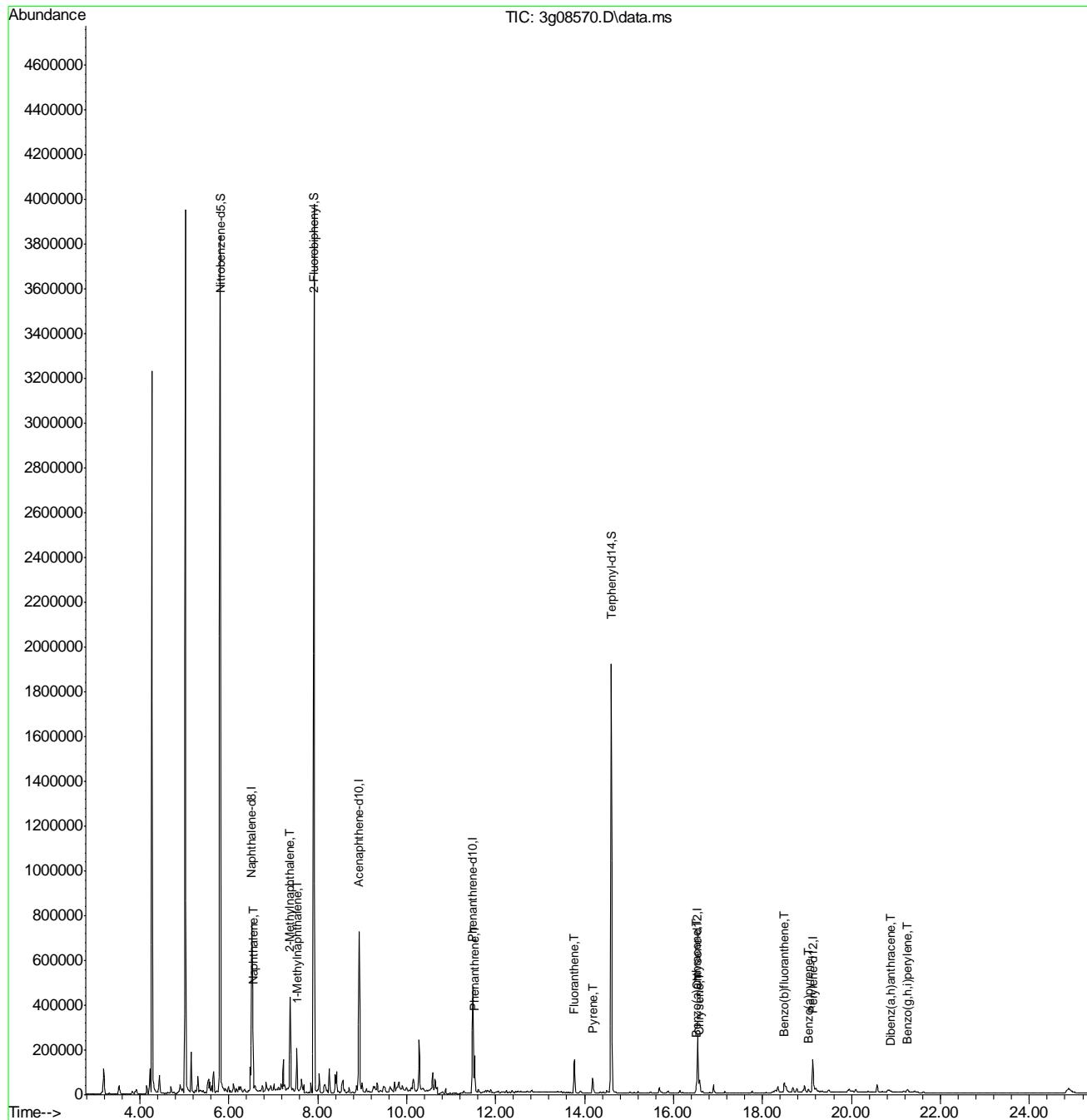
Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
<hr/>						
Internal Standards						
1) Naphthalene-d8	6.520	136	700667	4.00	ug/mL	-0.01
6) Acenaphthene-d10	8.933	164	379520	4.00	ug/mL	0.00
14) Phenanthrene-d10	11.485	188	507703	4.00	ug/mL	0.00
18) Chrysene-d12	16.547	240	290746	4.00	ug/mL	0.00
23) Perylene-d12	19.132	264	213019	4.00	ug/mL	0.00
<hr/>						
System Monitoring Compounds						
2) Nitrobenzene-d5	5.809	82	2372869	24.85	ug/mL	0.00
Spiked Amount 50.000	Range 25 - 135		Recovery	=	49.70%	
7) 2-Fluorobiphenyl	7.917	172	3373224	22.15	ug/mL	-0.01
Spiked Amount 50.000	Range 25 - 135		Recovery	=	44.30%	
20) Terphenyl-d14	14.603	244	2212264	35.20	ug/mL	0.00
Spiked Amount 50.000	Range 25 - 135		Recovery	=	70.40%	
<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	212235	0.90	ug/mL	97
8) 2-Methylnaphthalene	7.380	142	227418	1.65	ug/mL	98
9) 1-Methylnaphthalene	7.530	142	104712	0.79	ug/mL	97
10) Acenaphthylene	0.000		0	N.D.	d	
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	11.532	178	163892	0.93	ug/mL	99
16) Anthracene	0.000		0	N.D.	d	
17) Fluoranthene	13.772	202	109551	0.71	ug/mL	95
19) Pyrene	14.184	202	53513	0.41	ug/mL	89
21) Benzo(a)anthracene	16.514	228	44608	0.47	ug/mL	84
22) Chrysene	16.587	228	68812	0.69	ug/mL	92
24) Benzo(b)fluoranthene	18.491	252	52070m	0.68	ug/mL	
25) Benzo(k)fluoranthene	0.000		0	N.D.	d	
26) Benzo(a)pyrene	19.027	252	17574	0.24	ug/mL	92
27) Indeno(1,2,3-cd)pyrene	0.000		0	N.D.	d	
28) Dibenz(a,h)anthracene	20.888	278	6010m	0.16	ug/mL	
29) Benzo(g,h,i)perylene	21.267	276	17723	0.37	ug/mL	93
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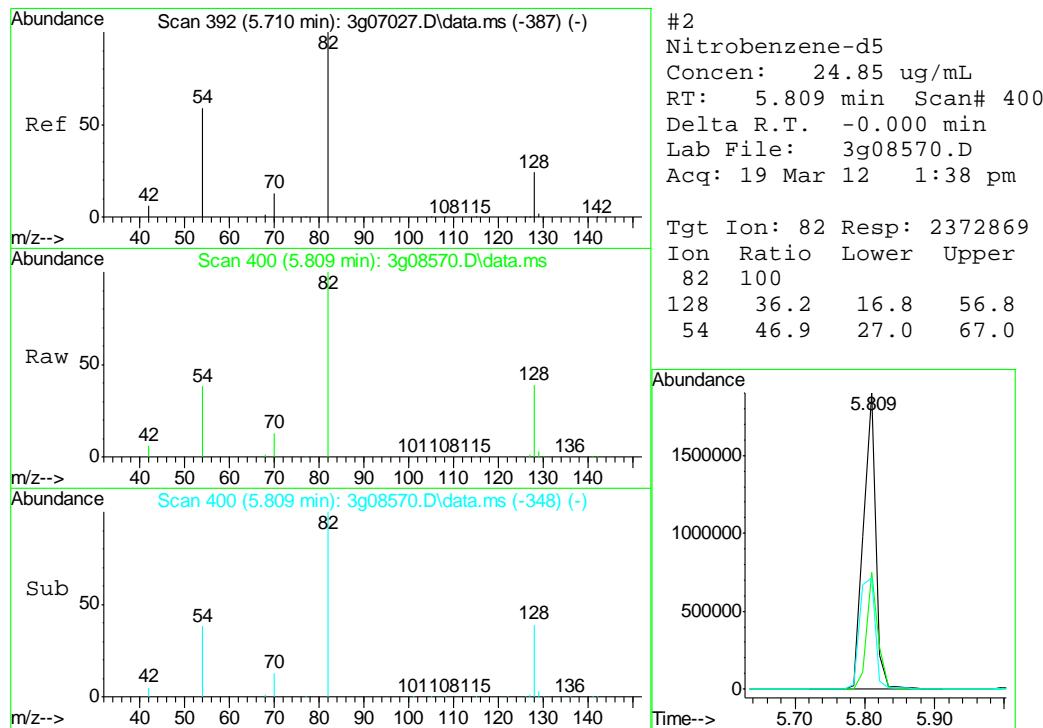
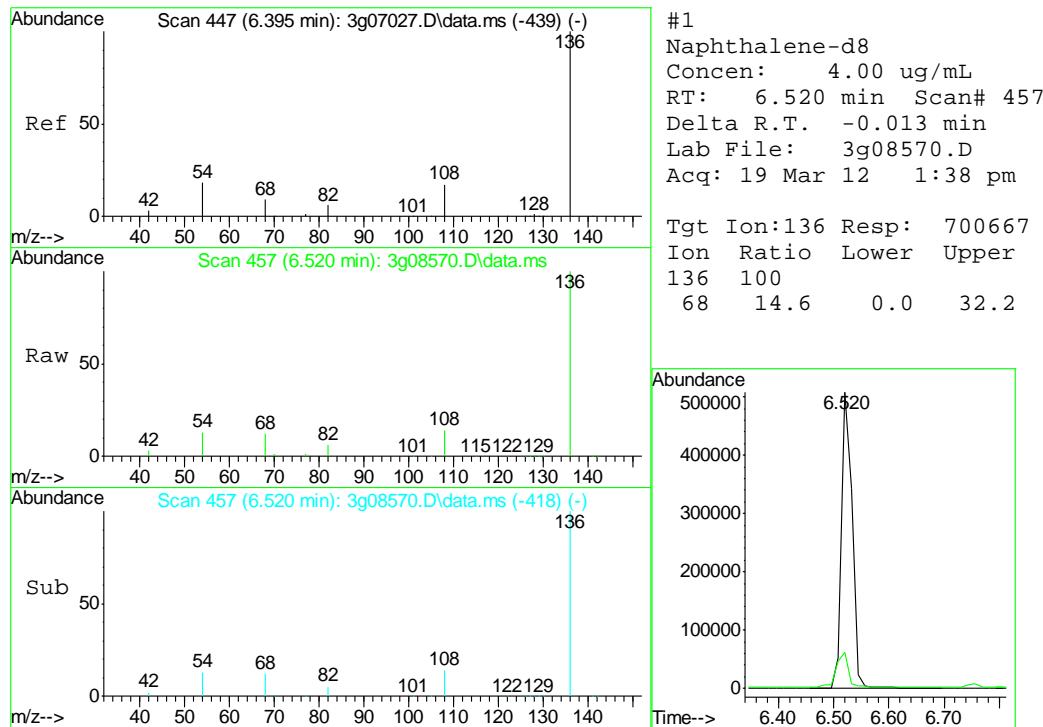
(#) = qualifier out of range (m) = manual integration (+) = signals summed

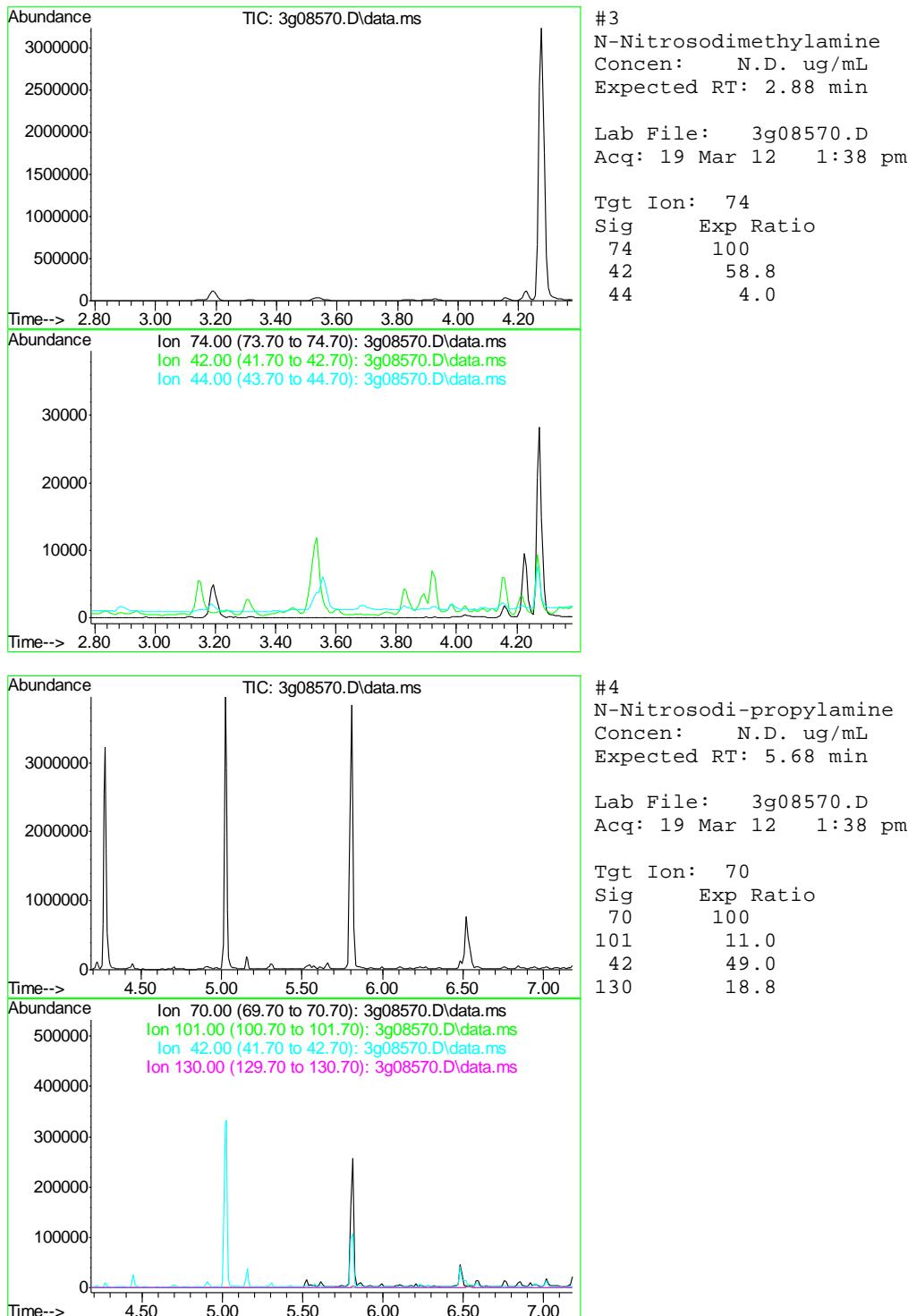
## Quantitation Report (QT Reviewed)

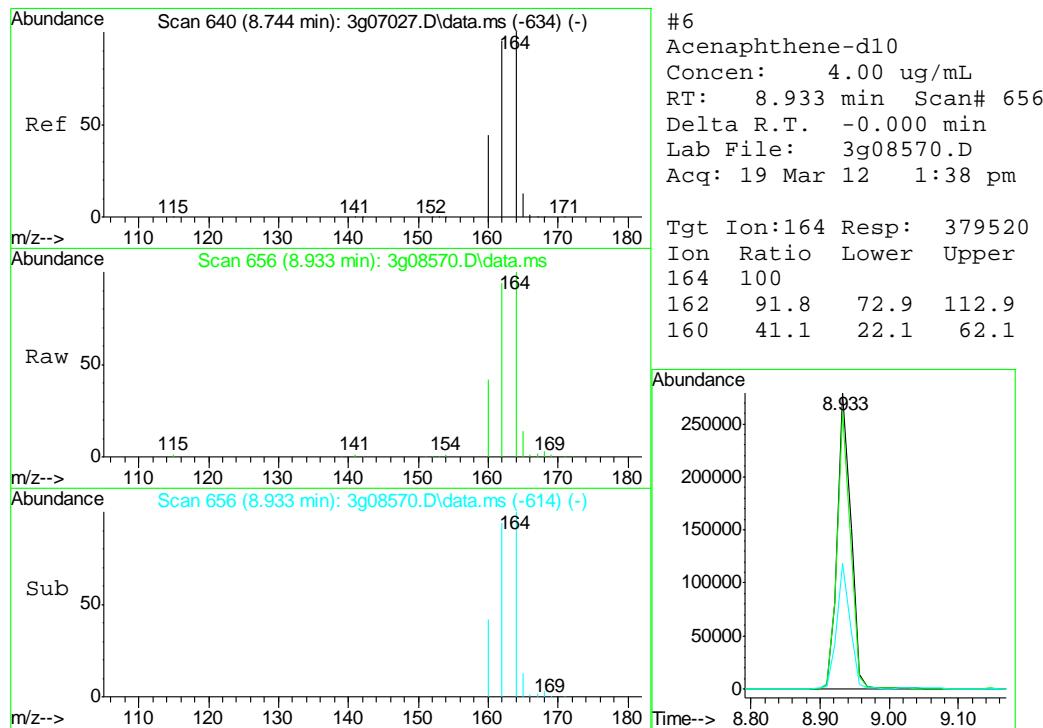
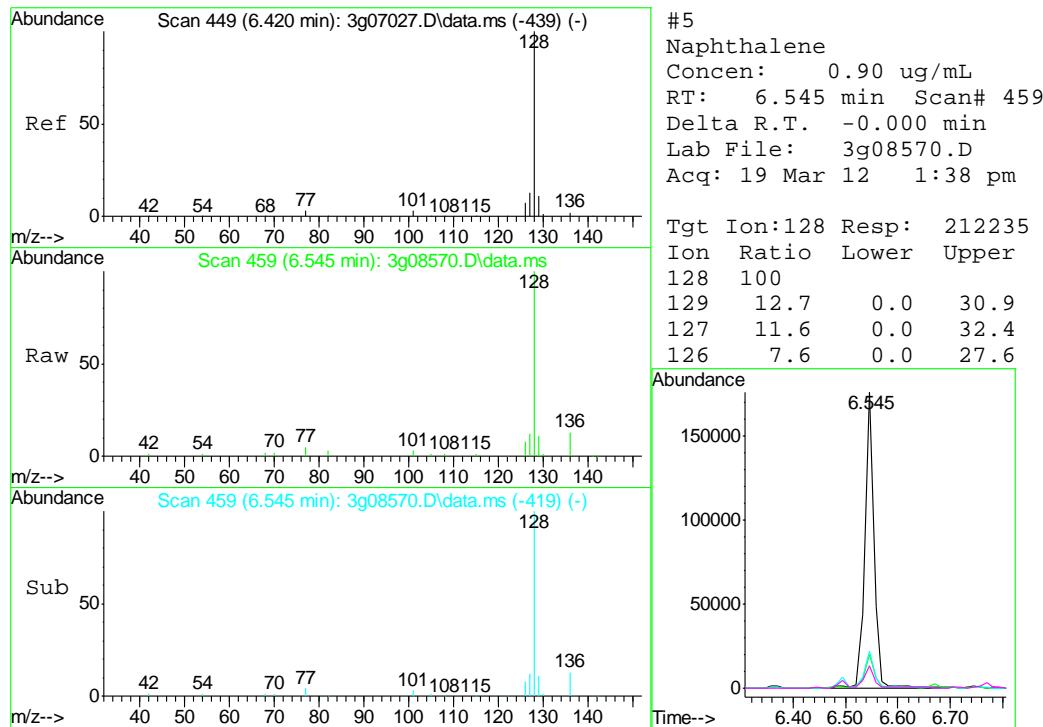
Data Path : C:\msdchem\1\DATA\031912\  
 Data File : 3g08570.D  
 Acq On : 19 Mar 2012 1:38 pm  
 Operator : DONC  
 Sample : D32747-1  
 Misc : OP5559,E3G352,30.02,,,1,1  
 ALS Vial : 6 Sample Multiplier: 1

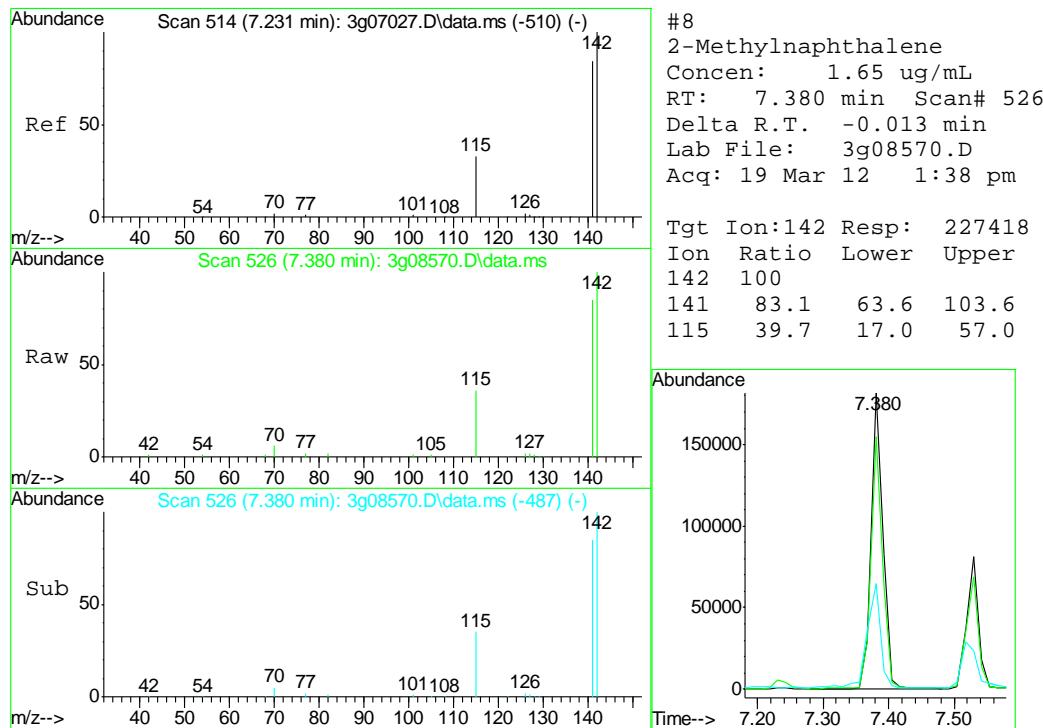
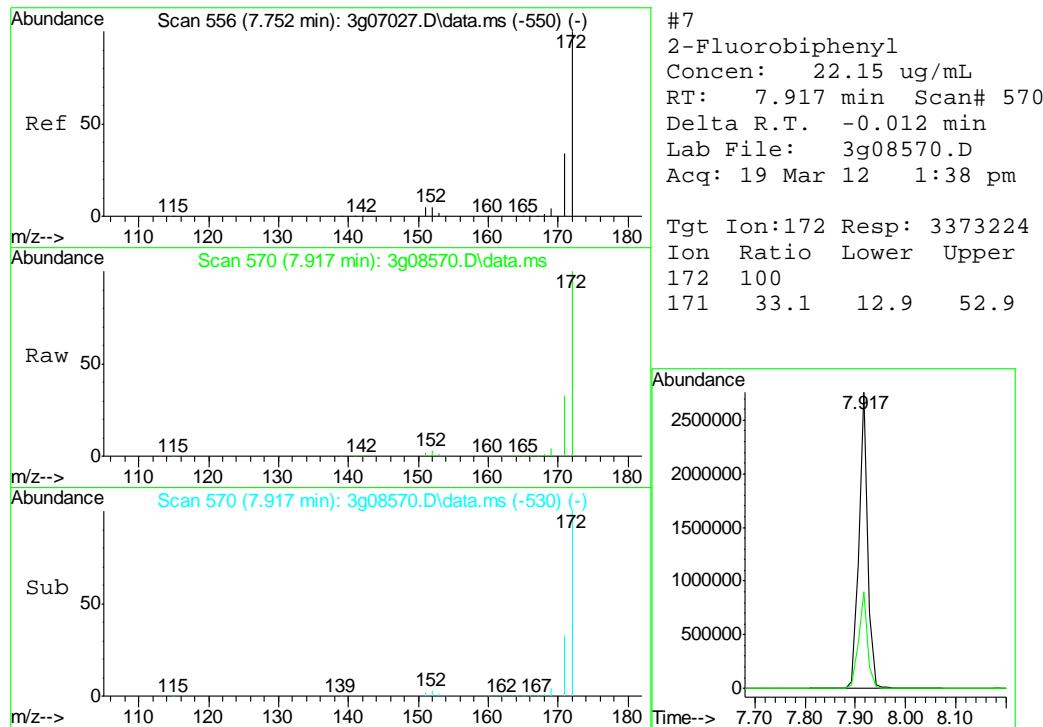
Quant Time: Mar 20 10:02:52 2012  
 Quant Method : C:\msdchem\1\METHODS\SIMPE3G344.M  
 Quant Title : PAHSIM BASE  
 QLast Update : Mon Mar 12 09:19:25 2012  
 Response via : Initial Calibration

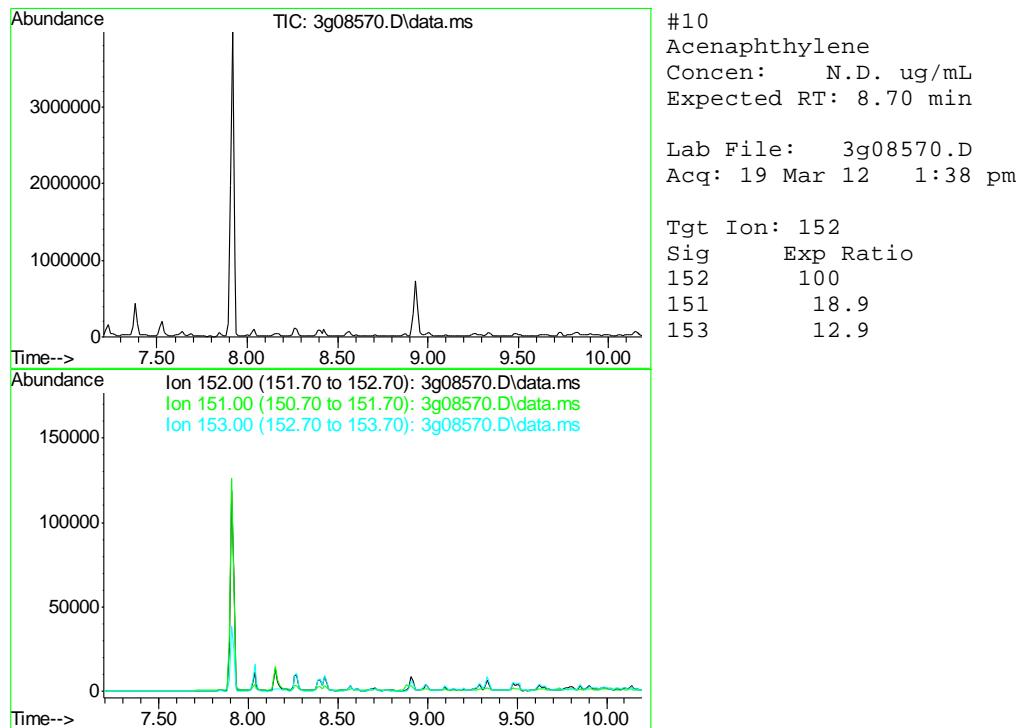
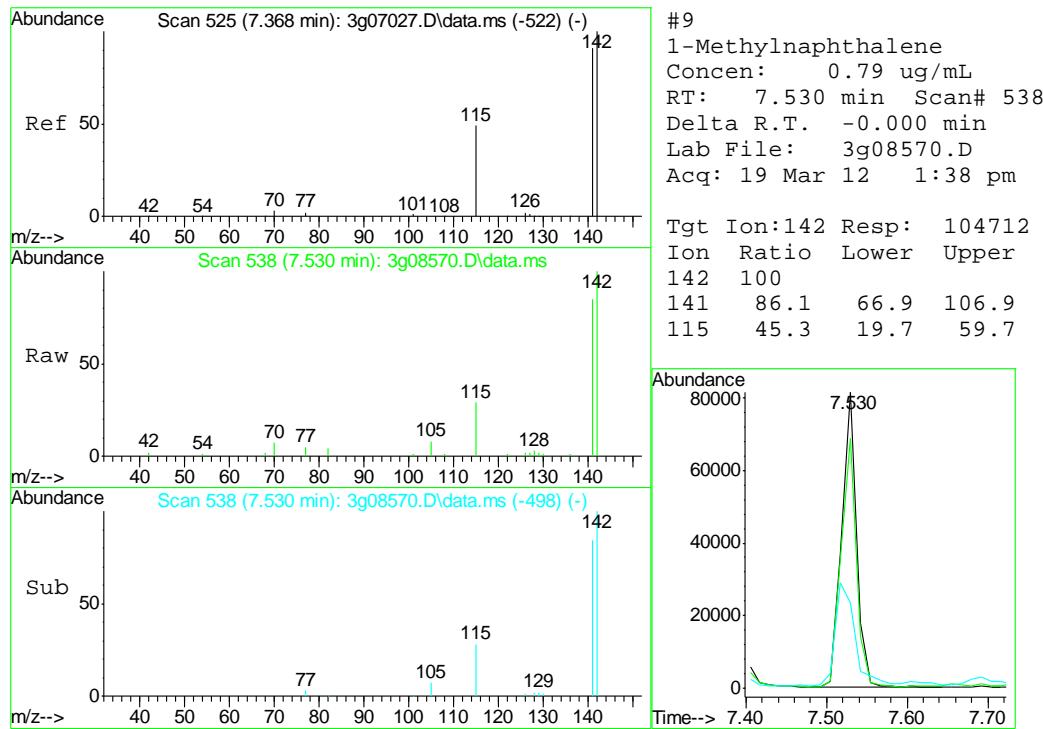


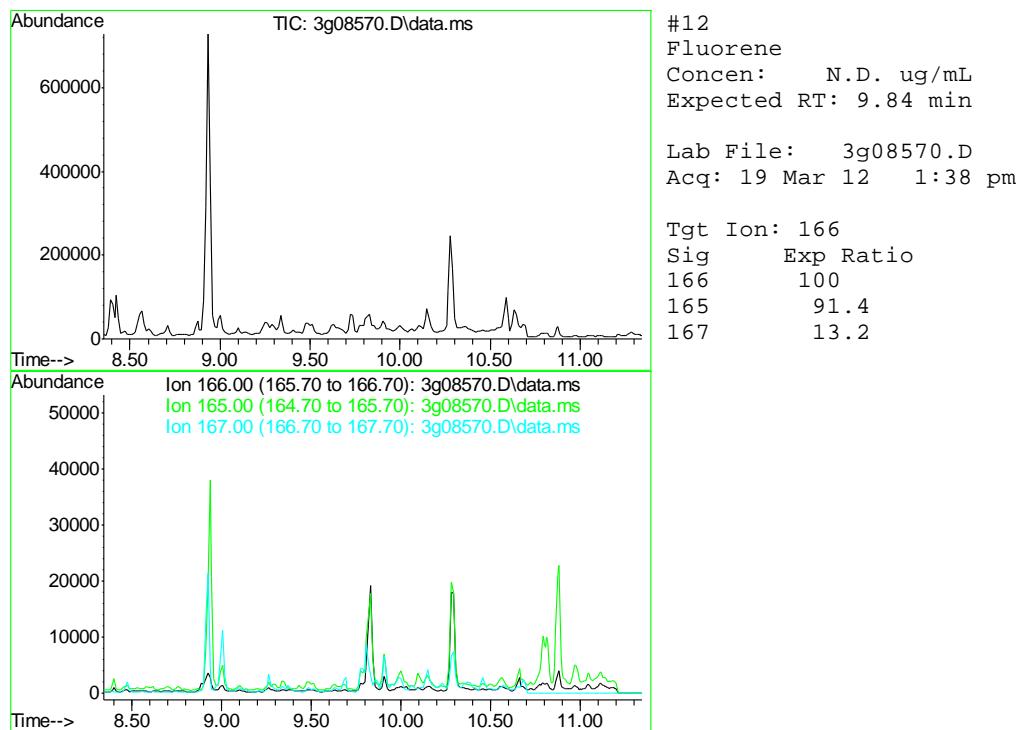
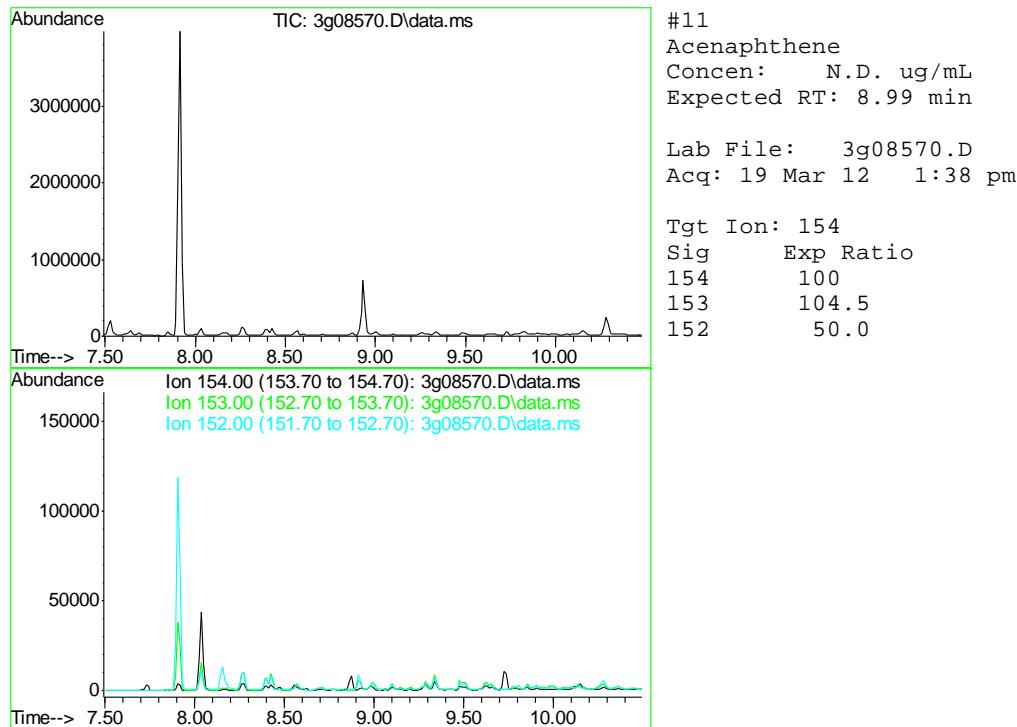


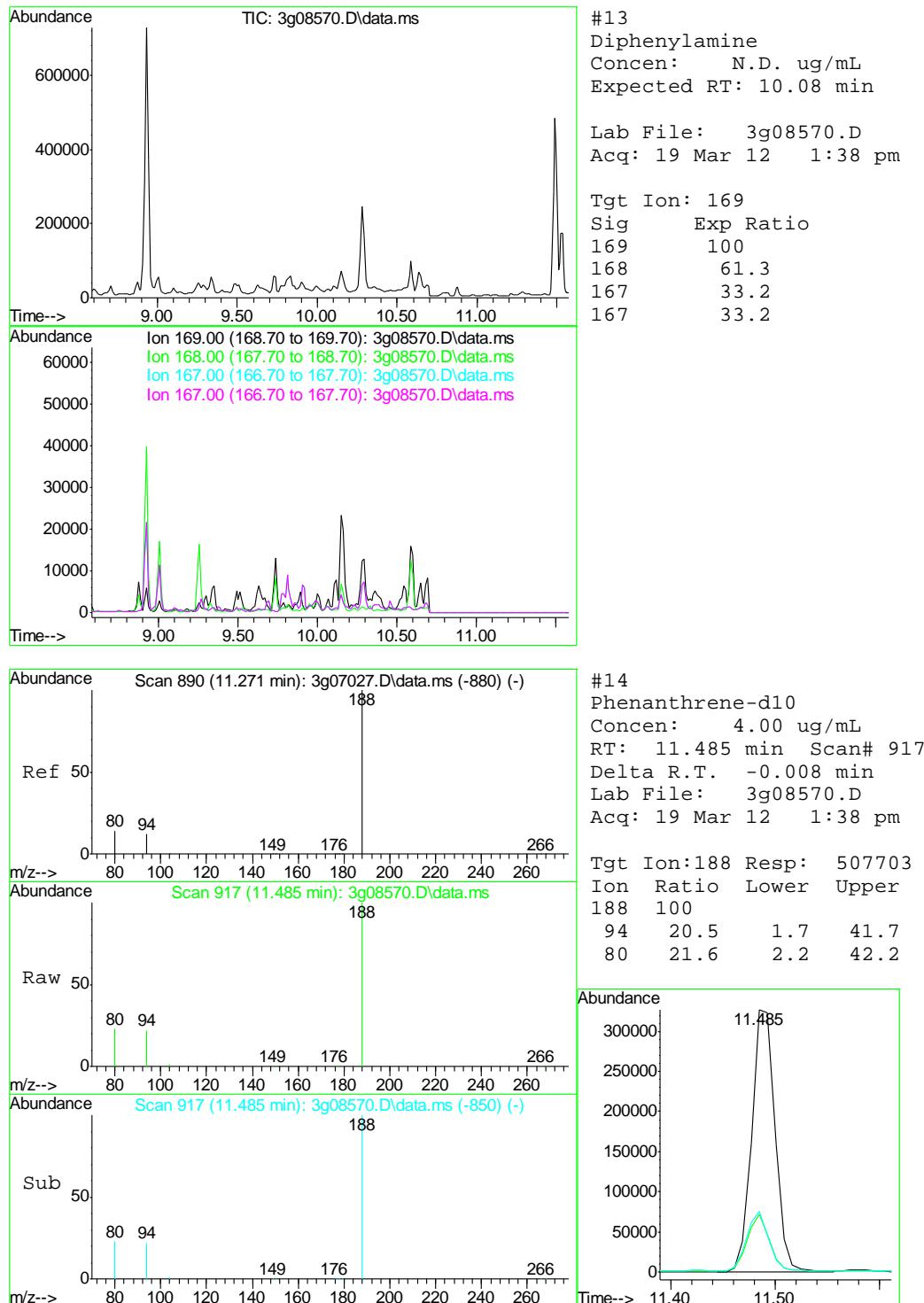


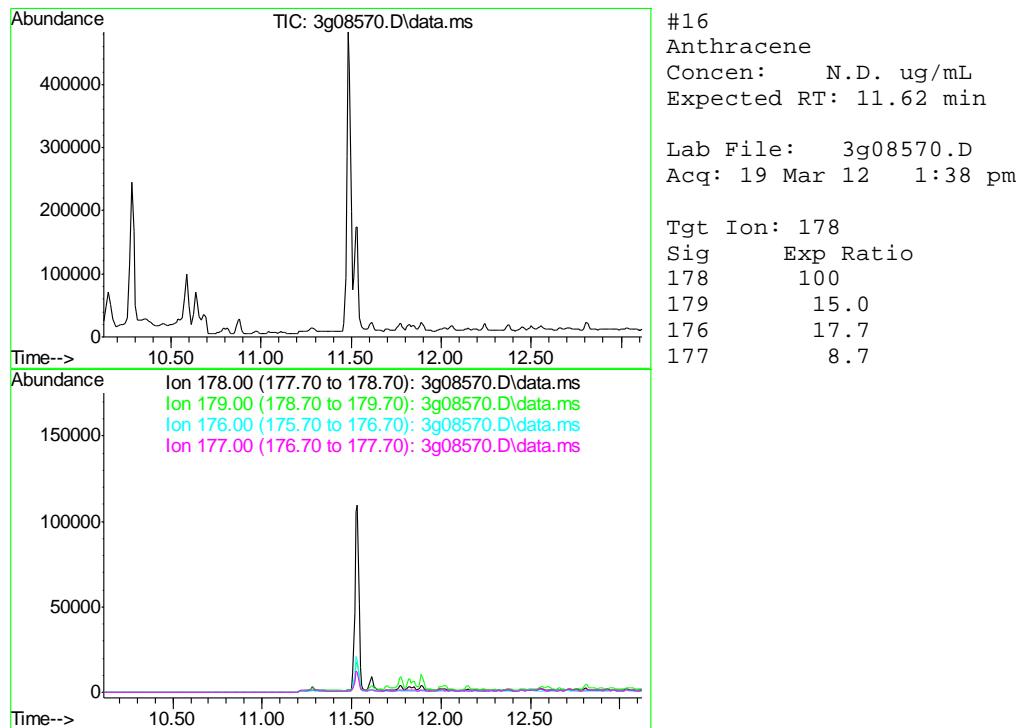
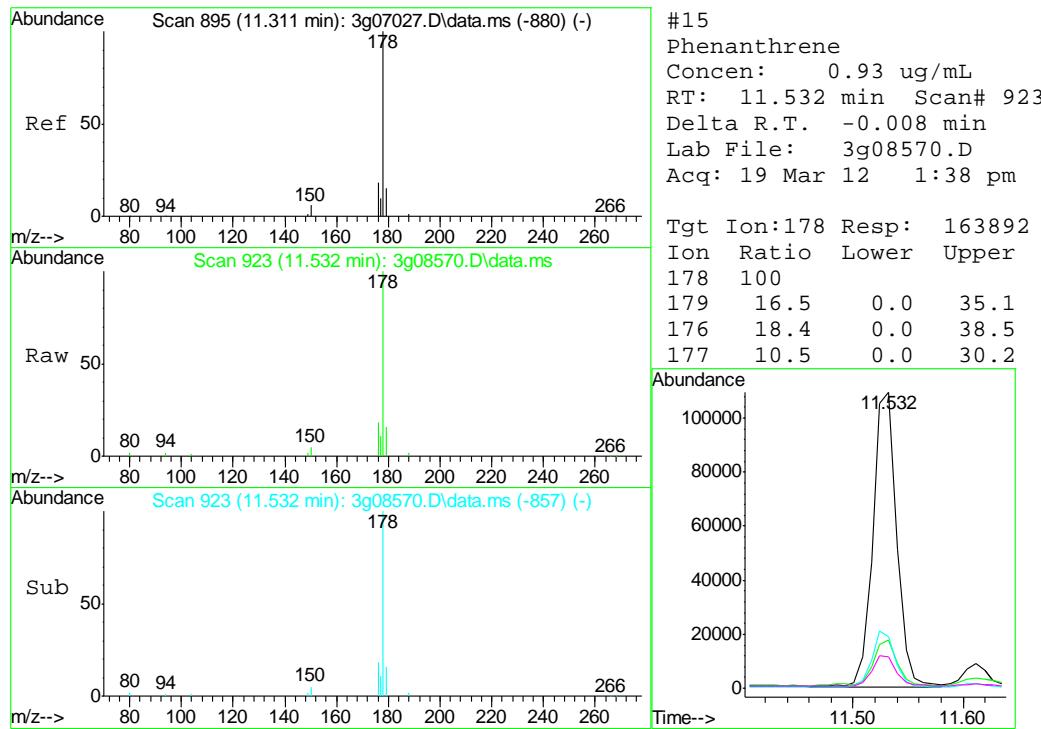


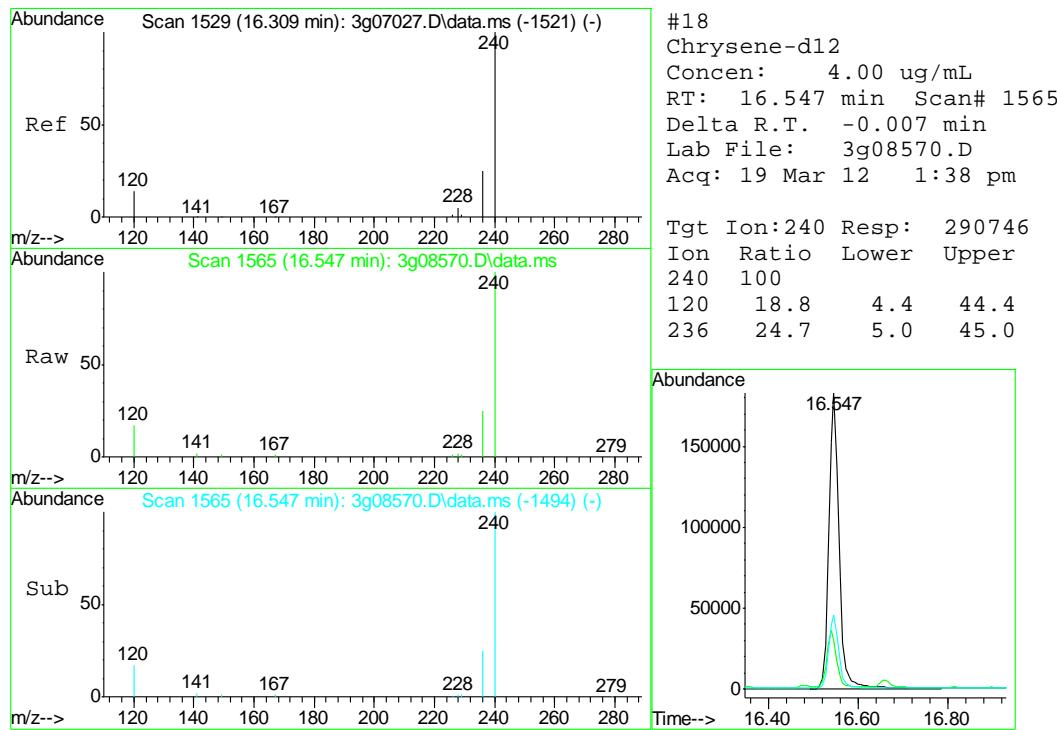
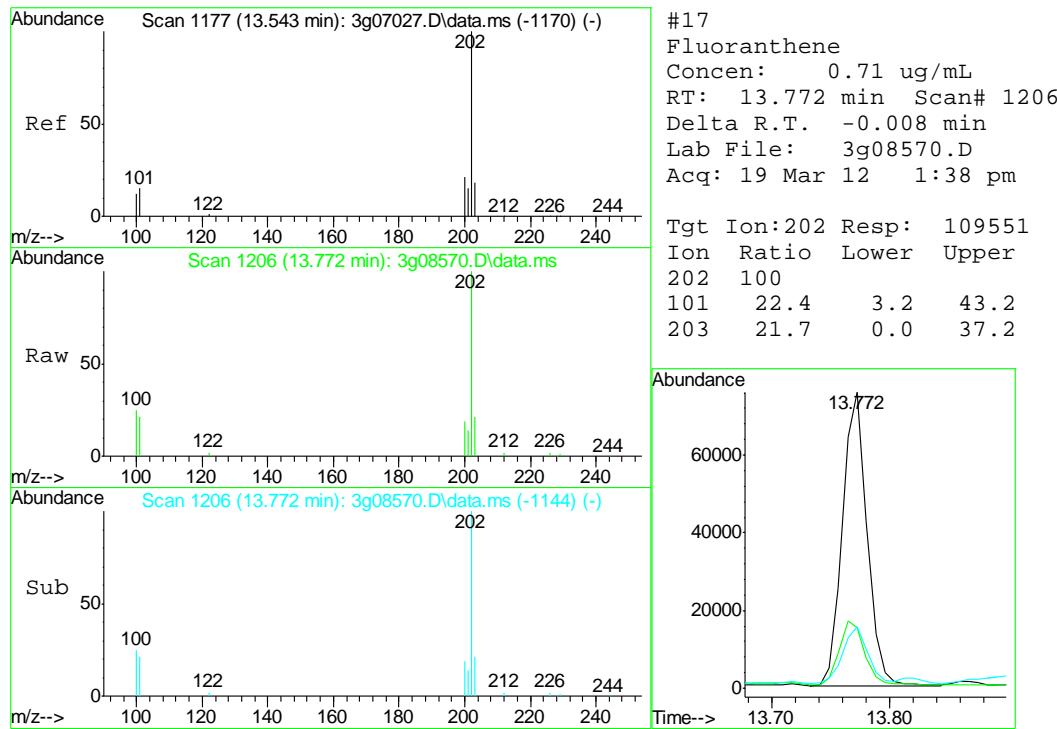


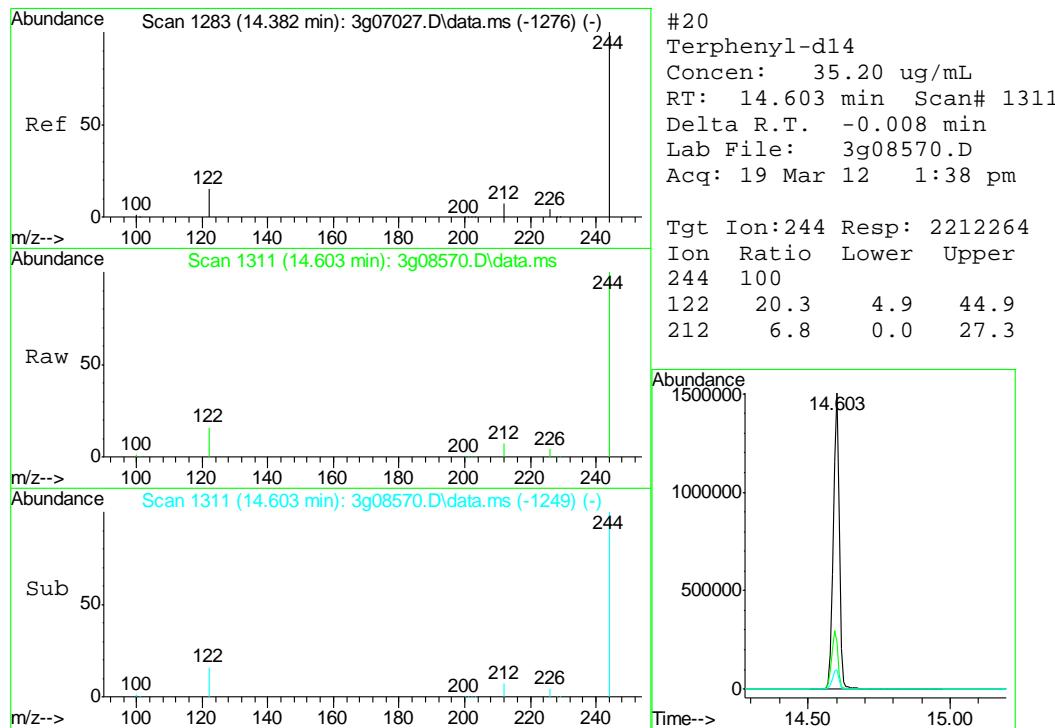
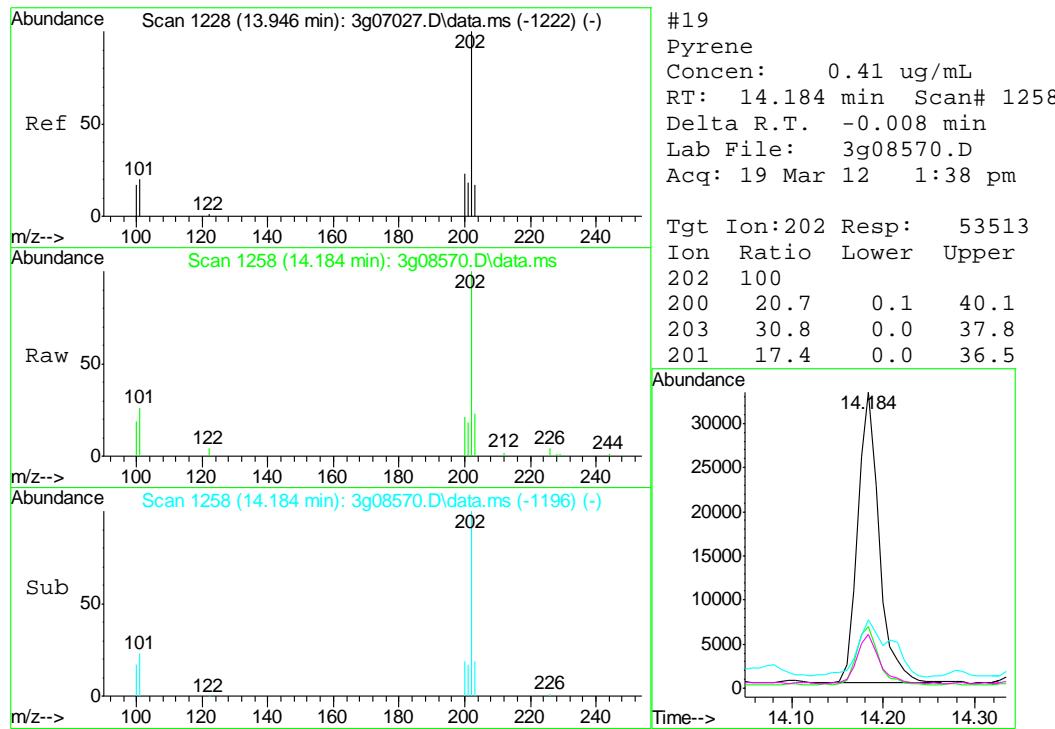


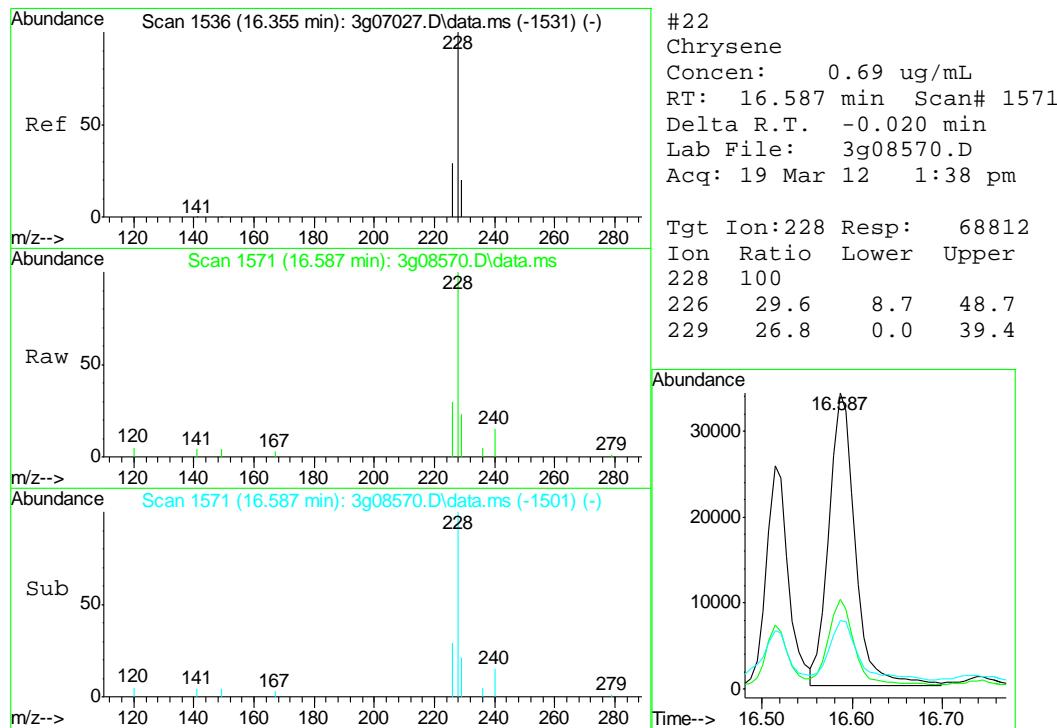
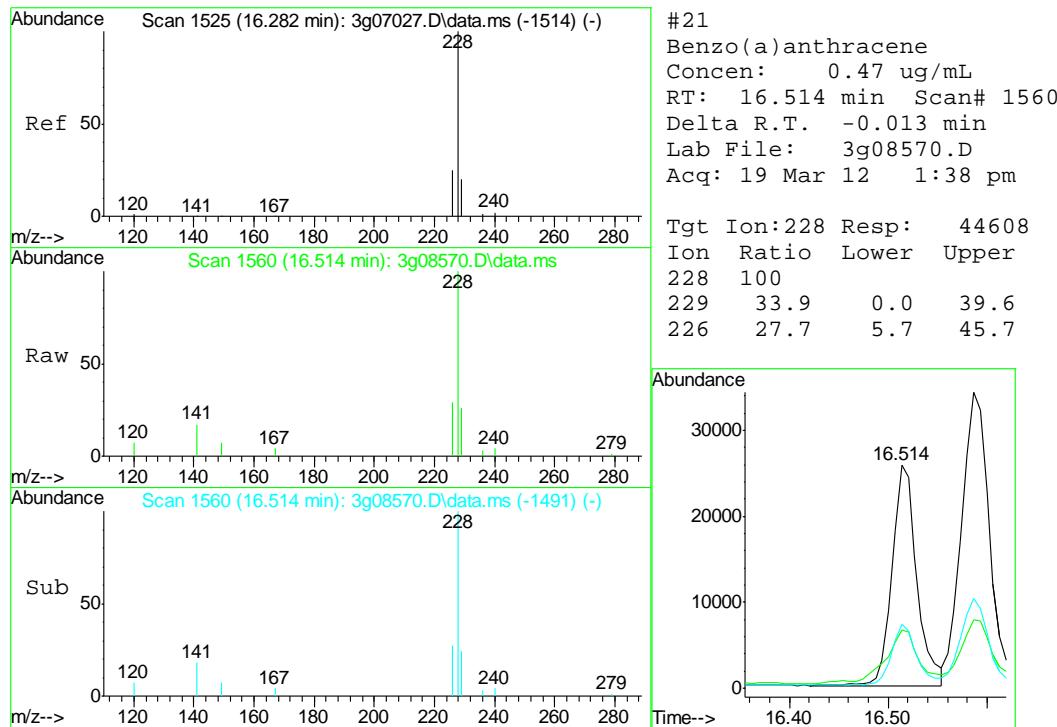


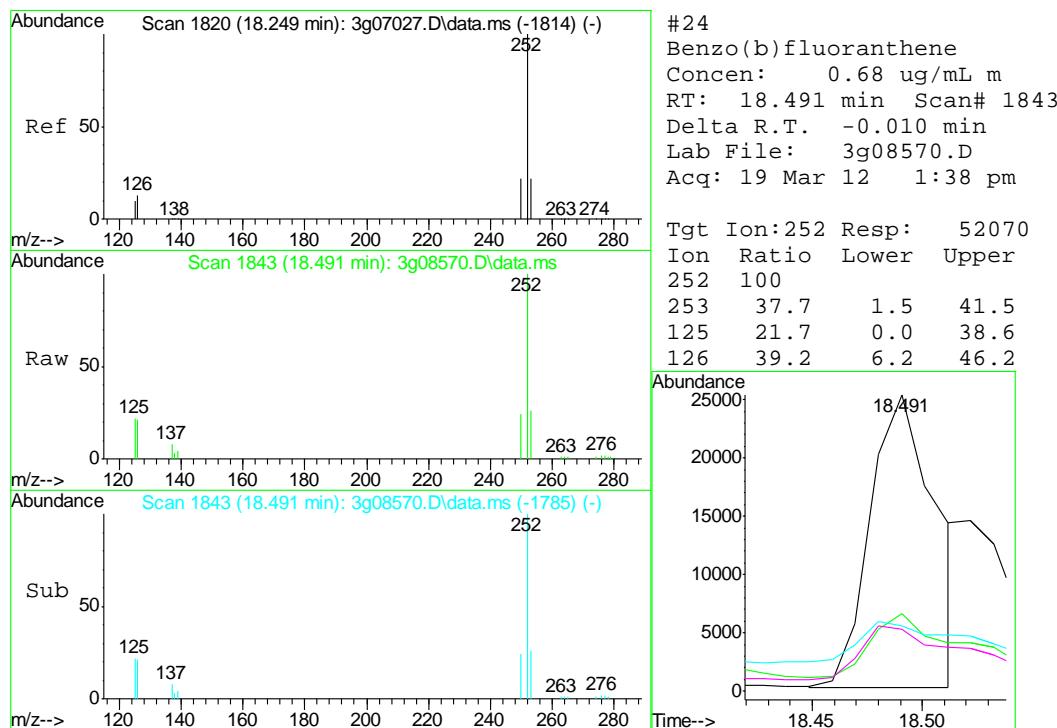
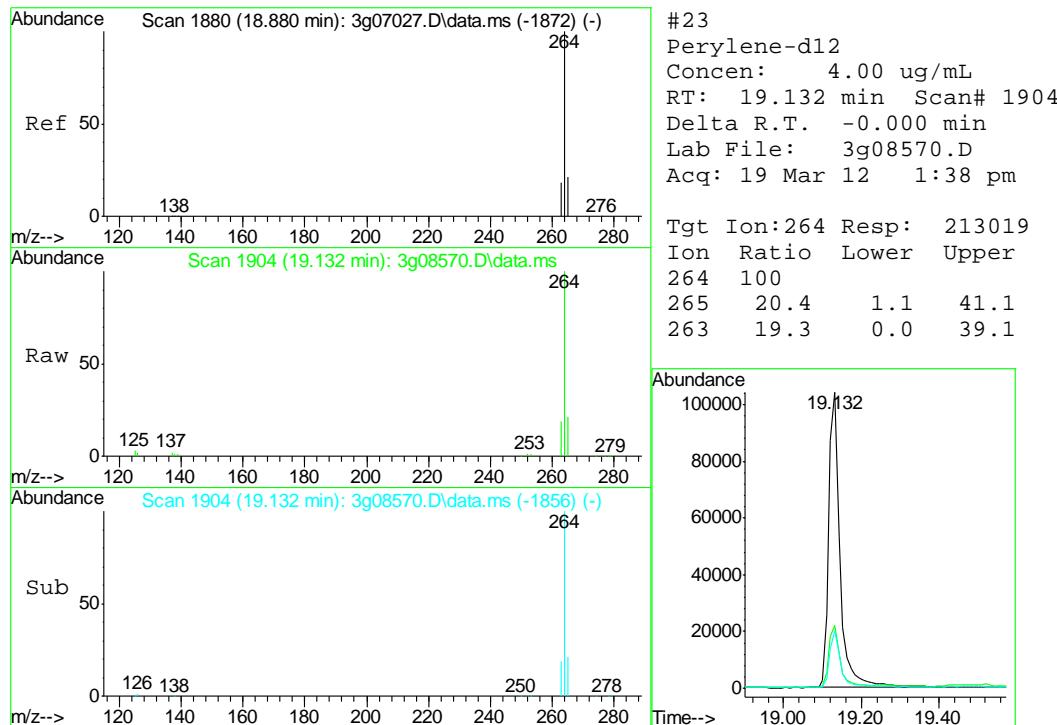


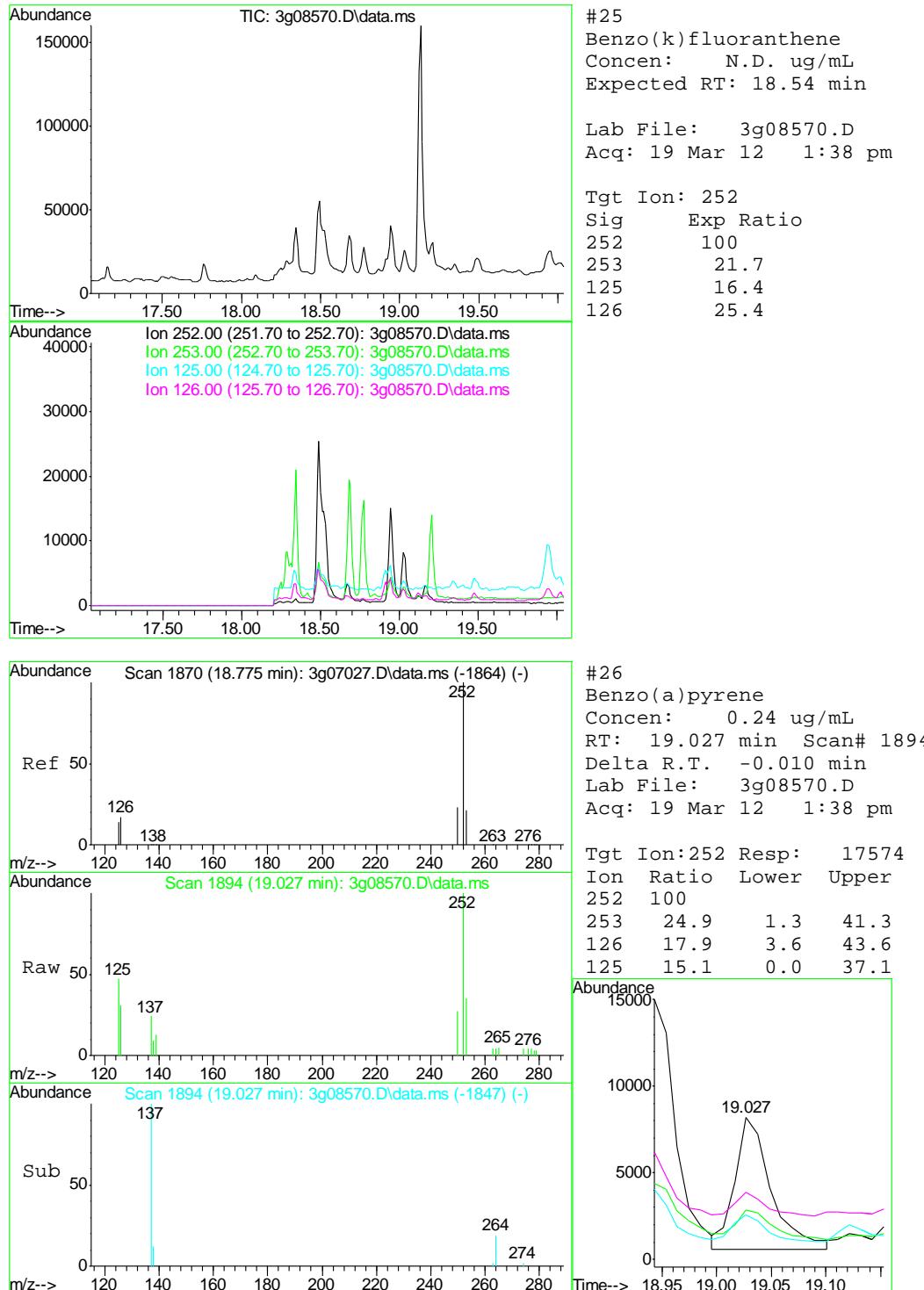


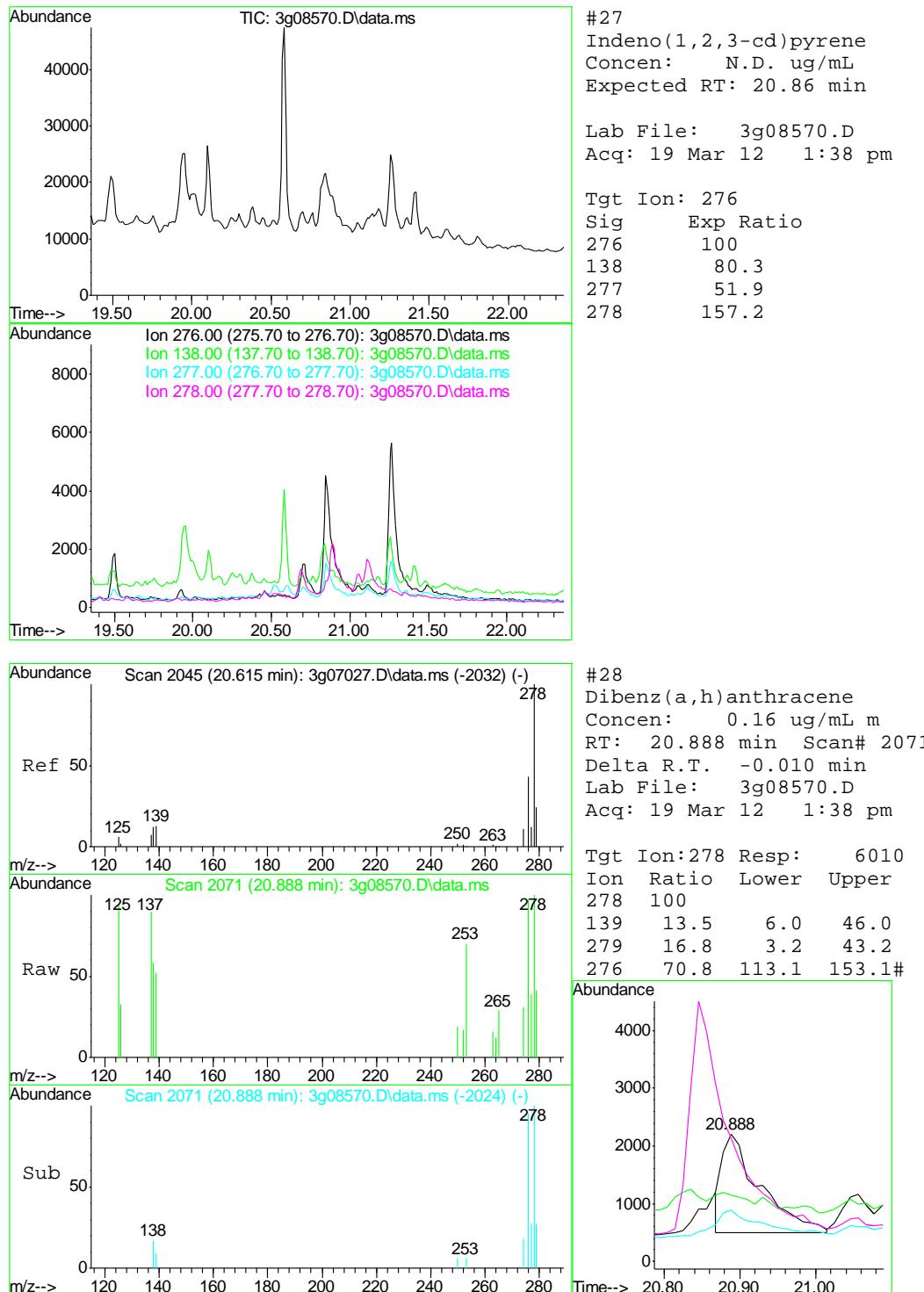


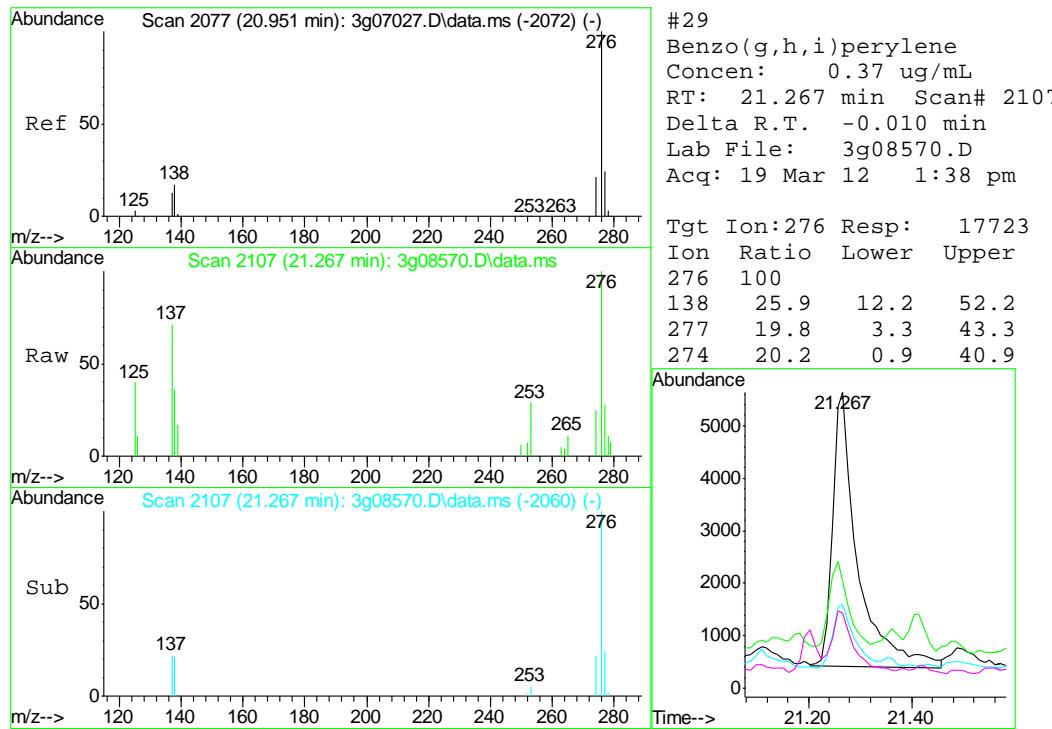












## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\031912\  
 Data File : 3g08568.D  
 Acq On : 19 Mar 2012 12:27 pm  
 Operator : DONC  
 Sample : OP5559-MB  
 Misc : OP5559,E3G352,30.00,,,1,1  
 ALS Vial : 4 Sample Multiplier: 1

Quant Time: Mar 20 09:59:44 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

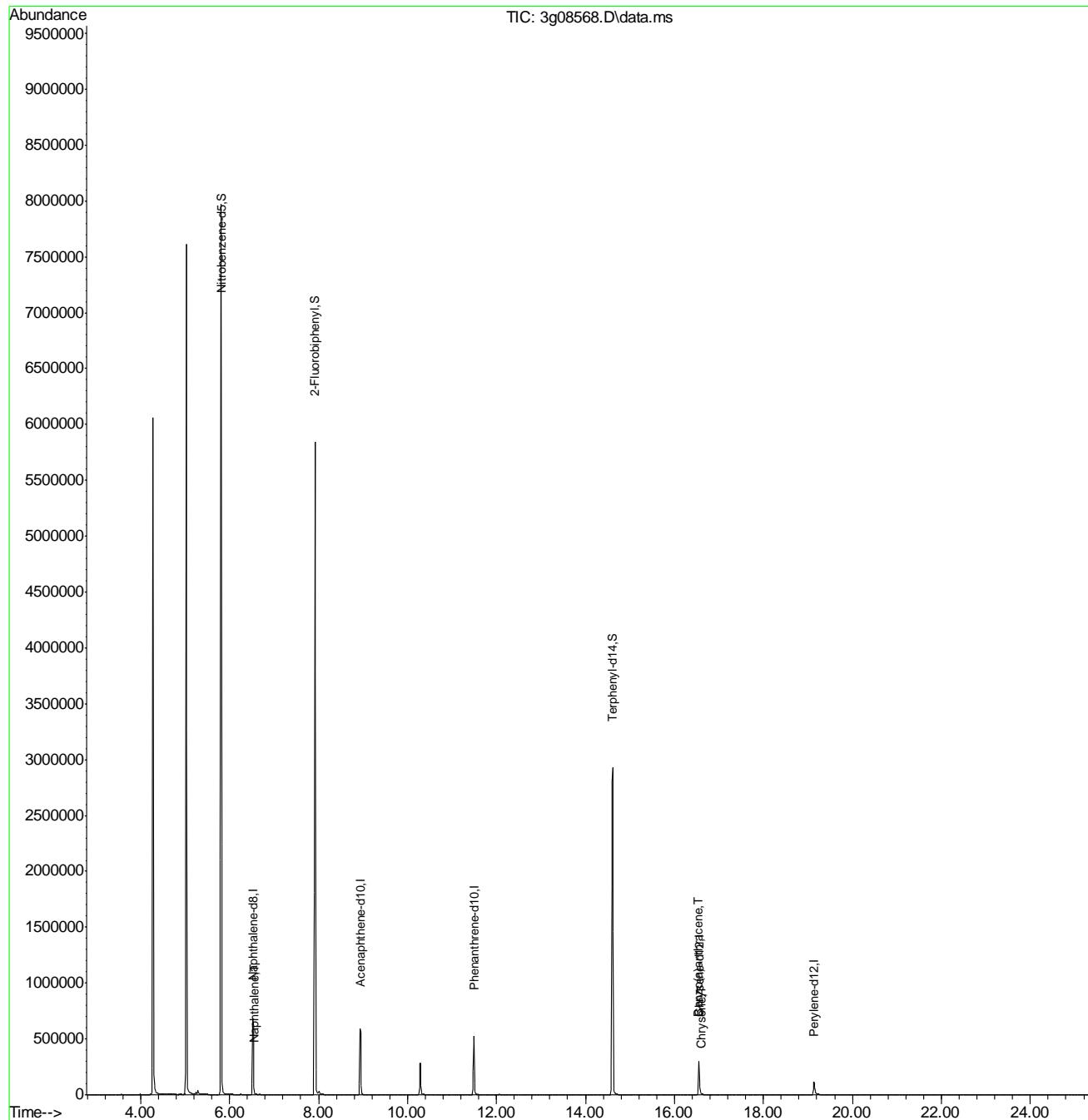
Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
<hr/>						
Internal Standards						
1) Naphthalene-d8	6.532	136	768016	4.00	ug/mL	0.00
6) Acenaphthene-d10	8.945	164	394036	4.00	ug/mL	0.01
14) Phenanthrene-d10	11.493	188	564653	4.00	ug/mL	0.00
18) Chrysene-d12	16.553	240	381549	4.00	ug/mL	0.00
23) Perylene-d12	19.132	264	207065	4.00	ug/mL	0.00
<hr/>						
System Monitoring Compounds						
2) Nitrobenzene-d5	5.809	82	4649774	44.43	ug/mL	0.00
Spiked Amount 50.000	Range 25 - 135		Recovery	=	88.86%	
7) 2-Fluorobiphenyl	7.917	172	5637717	35.66	ug/mL	-0.01
Spiked Amount 50.000	Range 25 - 135		Recovery	=	71.32%	
20) Terphenyl-d14	14.611	244	3706451	44.94	ug/mL	0.00
Spiked Amount 50.000	Range 25 - 135		Recovery	=	89.88%	
<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	1341	0.01	ug/mL#	73
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.	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	16.547	228	1429	0.01	ug/mL	68
22) Chrysene	16.600	228	865	0.01	ug/mL#	71
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	

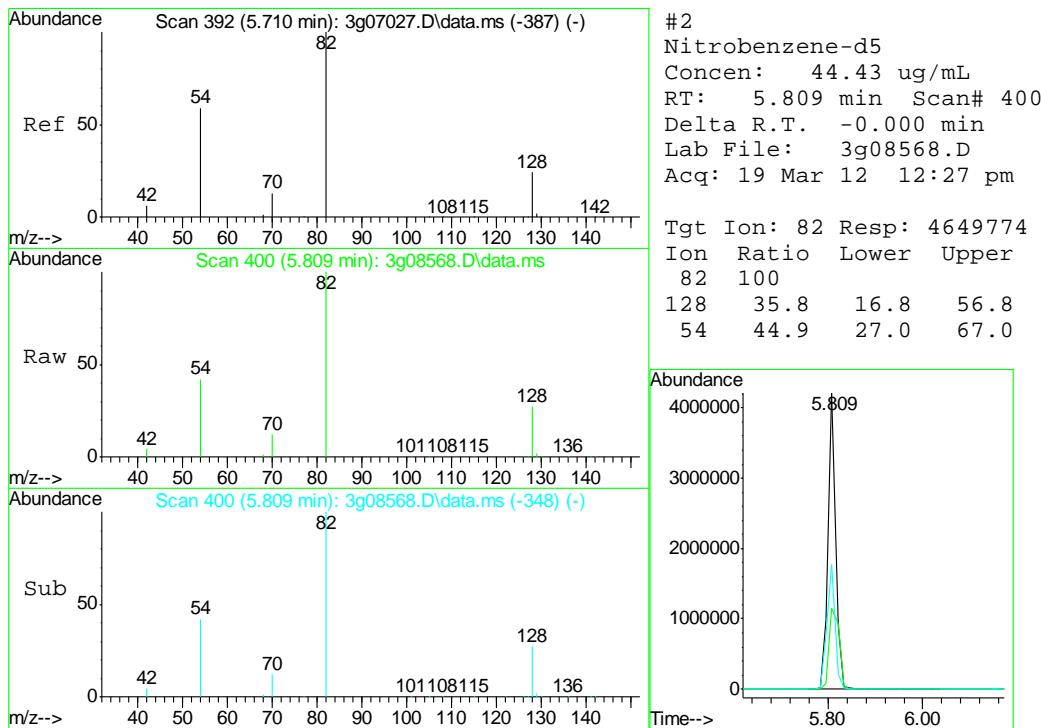
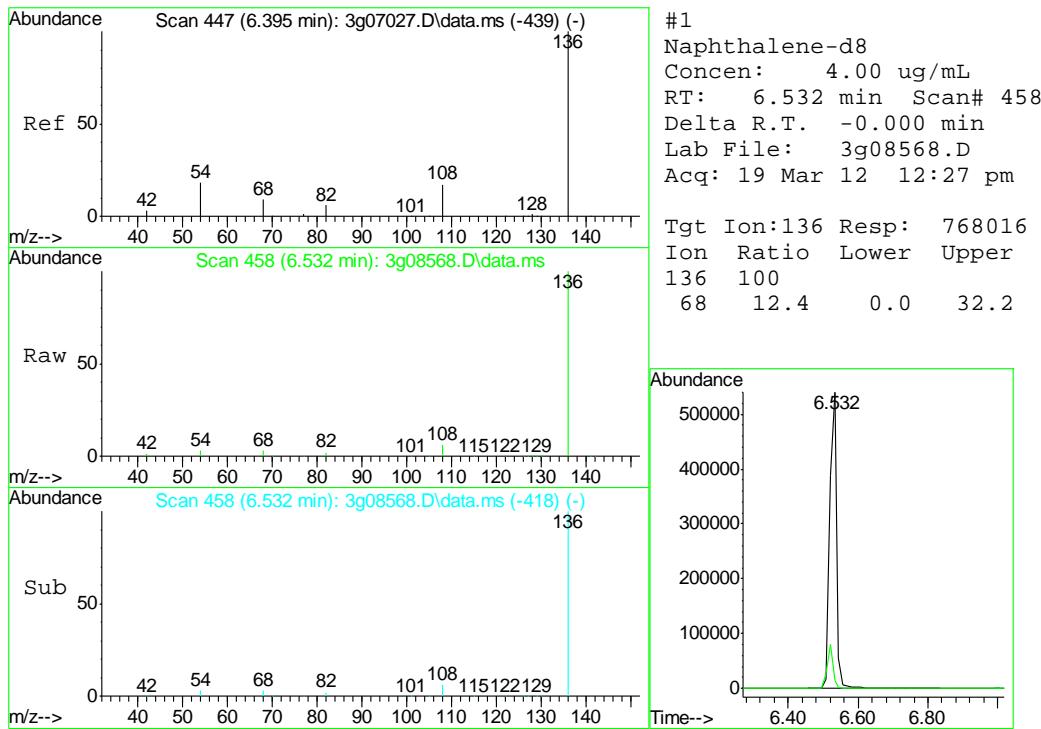
(#) = qualifier out of range (m) = manual integration (+) = signals summed

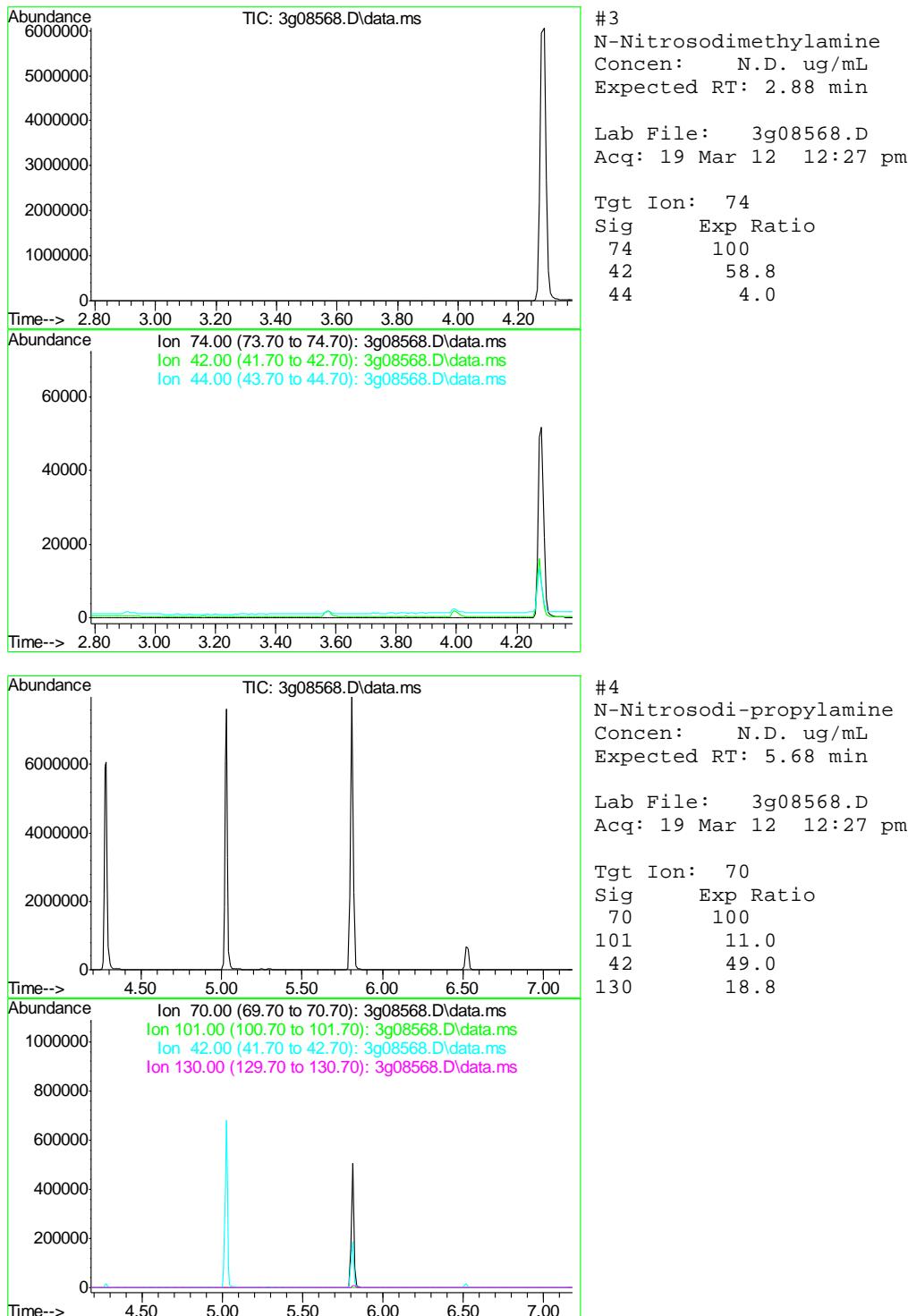
## Quantitation Report (QT Reviewed)

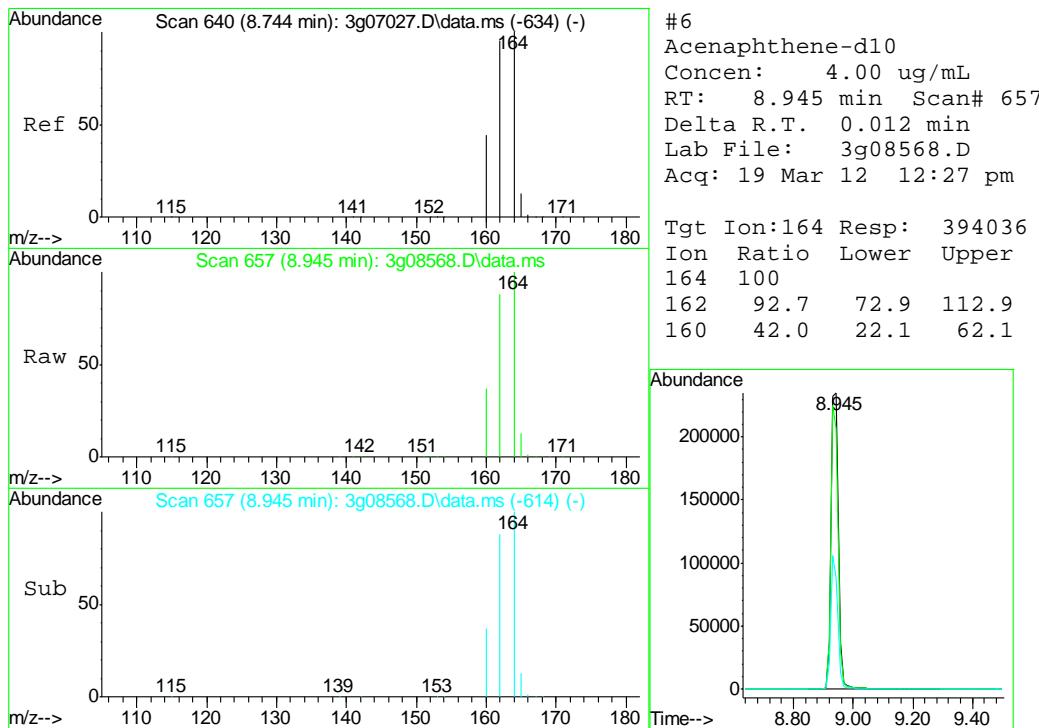
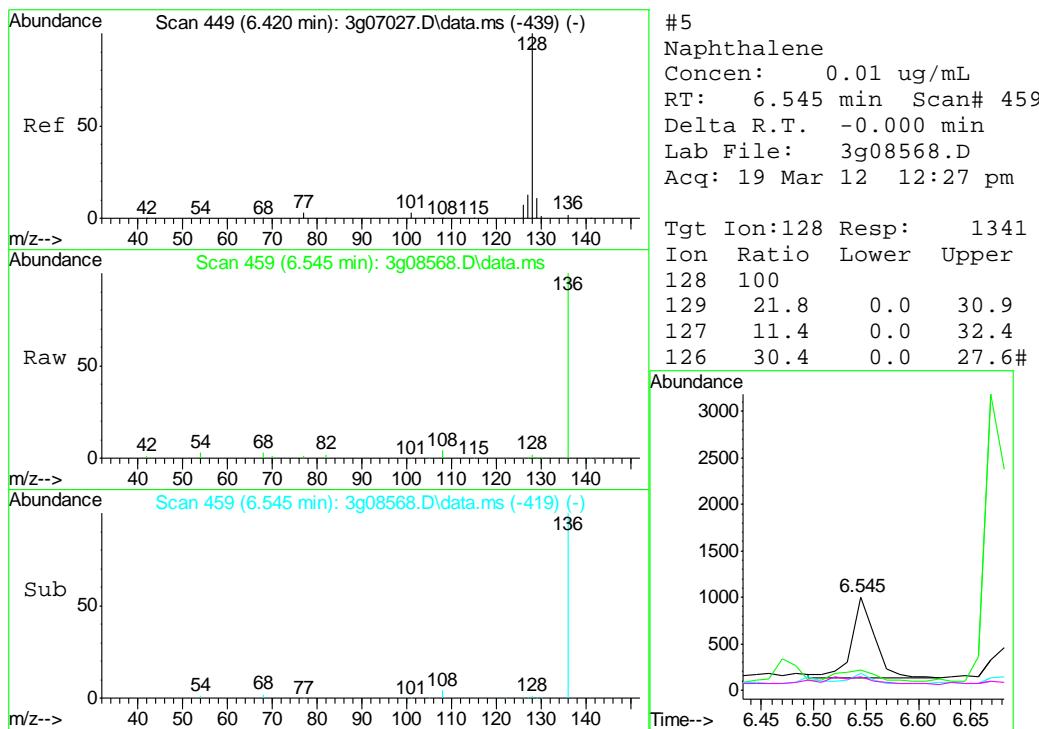
Data Path : C:\msdchem\1\DATA\031912\  
 Data File : 3g08568.D  
 Acq On : 19 Mar 2012 12:27 pm  
 Operator : DONC  
 Sample : OP5559-MB  
 Misc : OP5559,E3G352,30.00,,,1,1  
 ALS Vial : 4 Sample Multiplier: 1

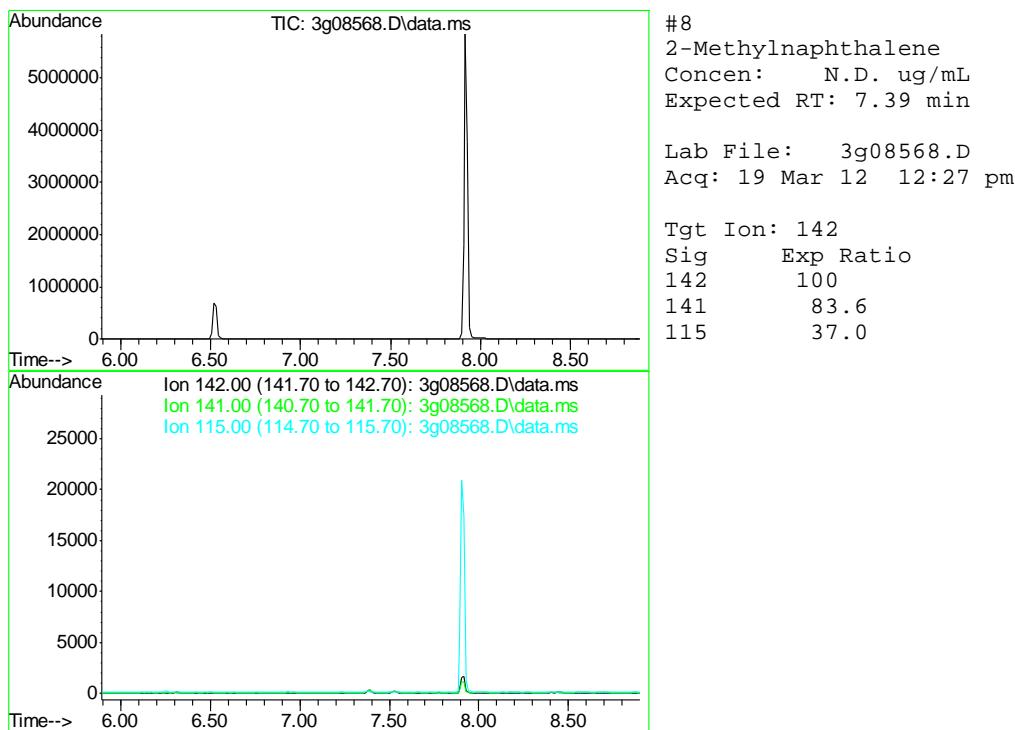
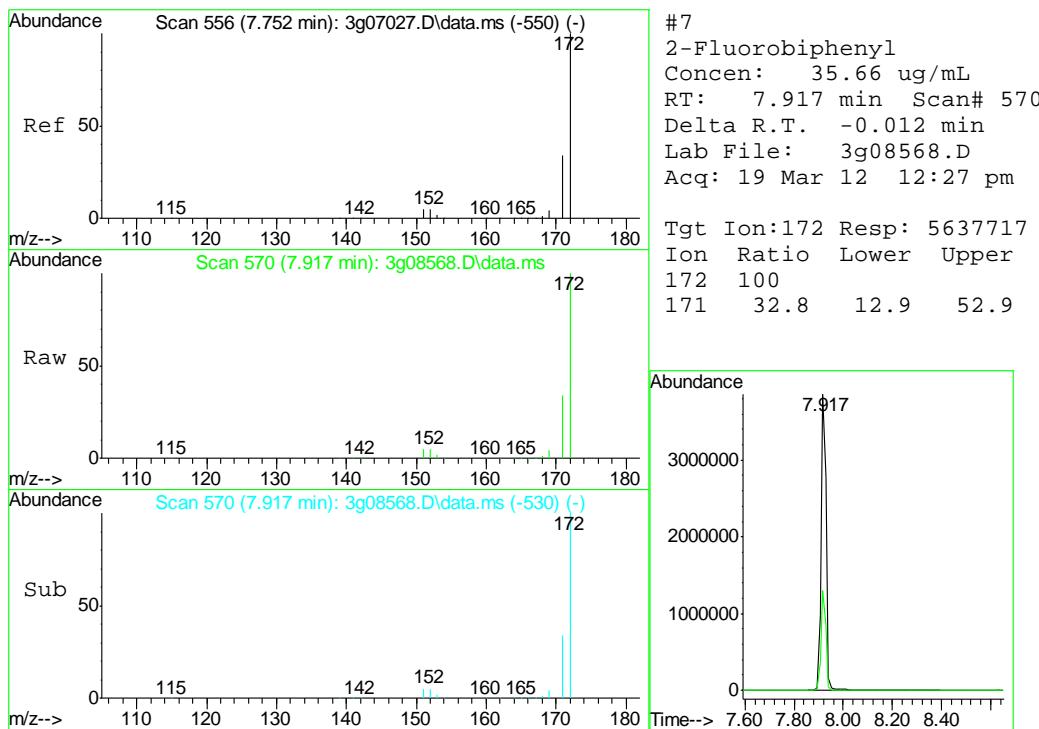
Quant Time: Mar 20 09:59:44 2012  
 Quant Method : C:\msdchem\1\METHODS\SIMPE3G344.M  
 Quant Title : PAHSIM BASE  
 QLast Update : Mon Mar 12 09:19:25 2012  
 Response via : Initial Calibration

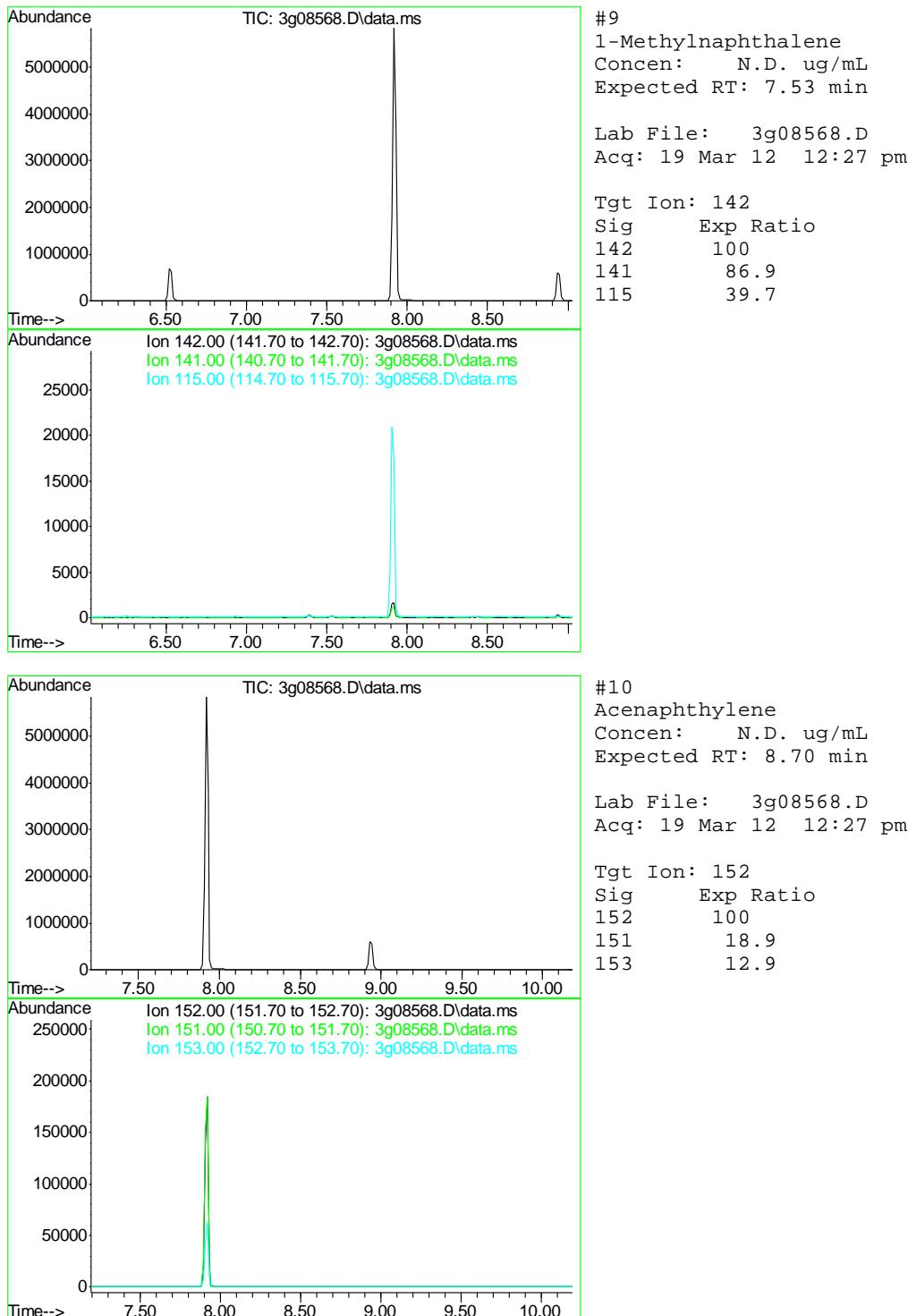


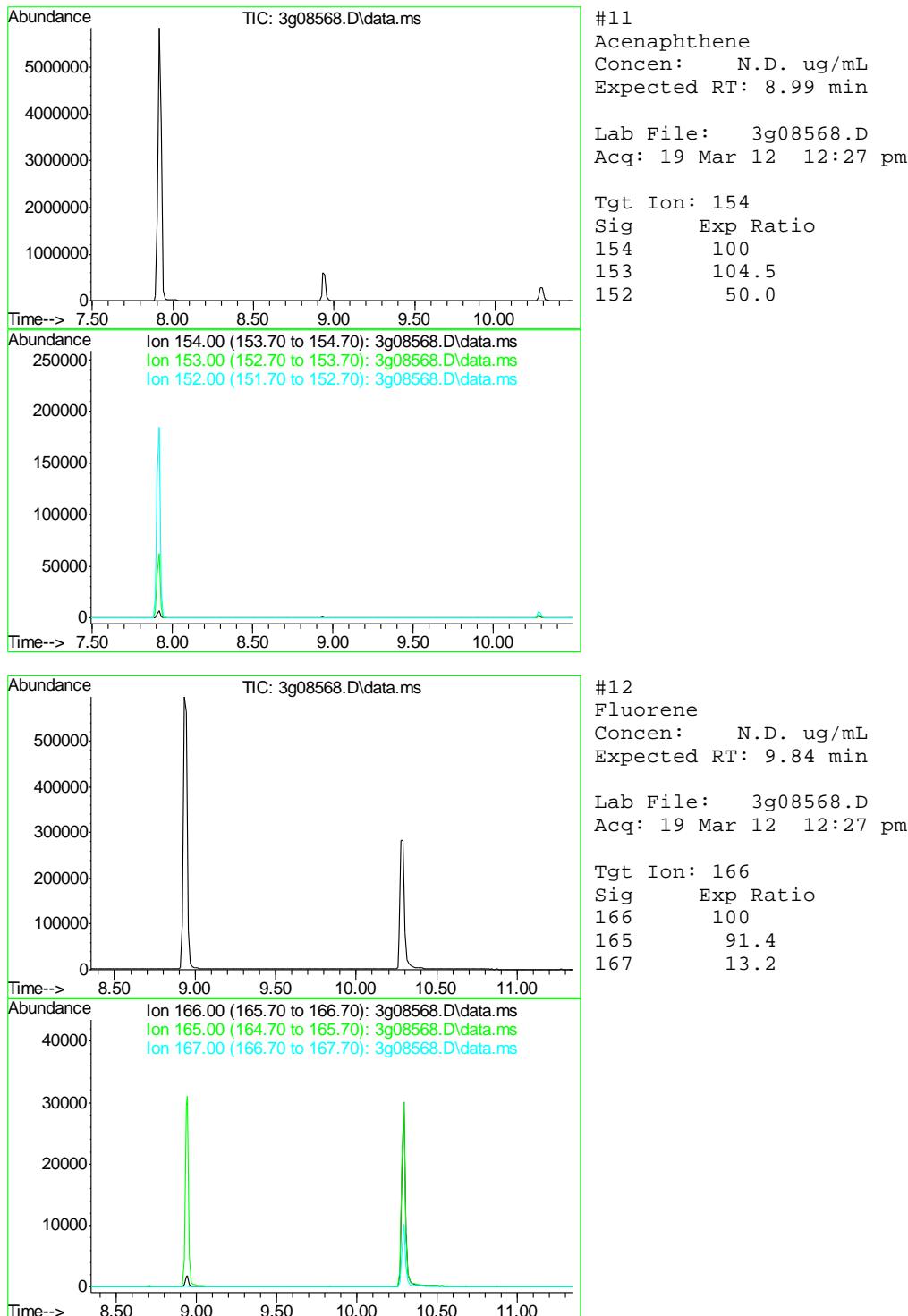


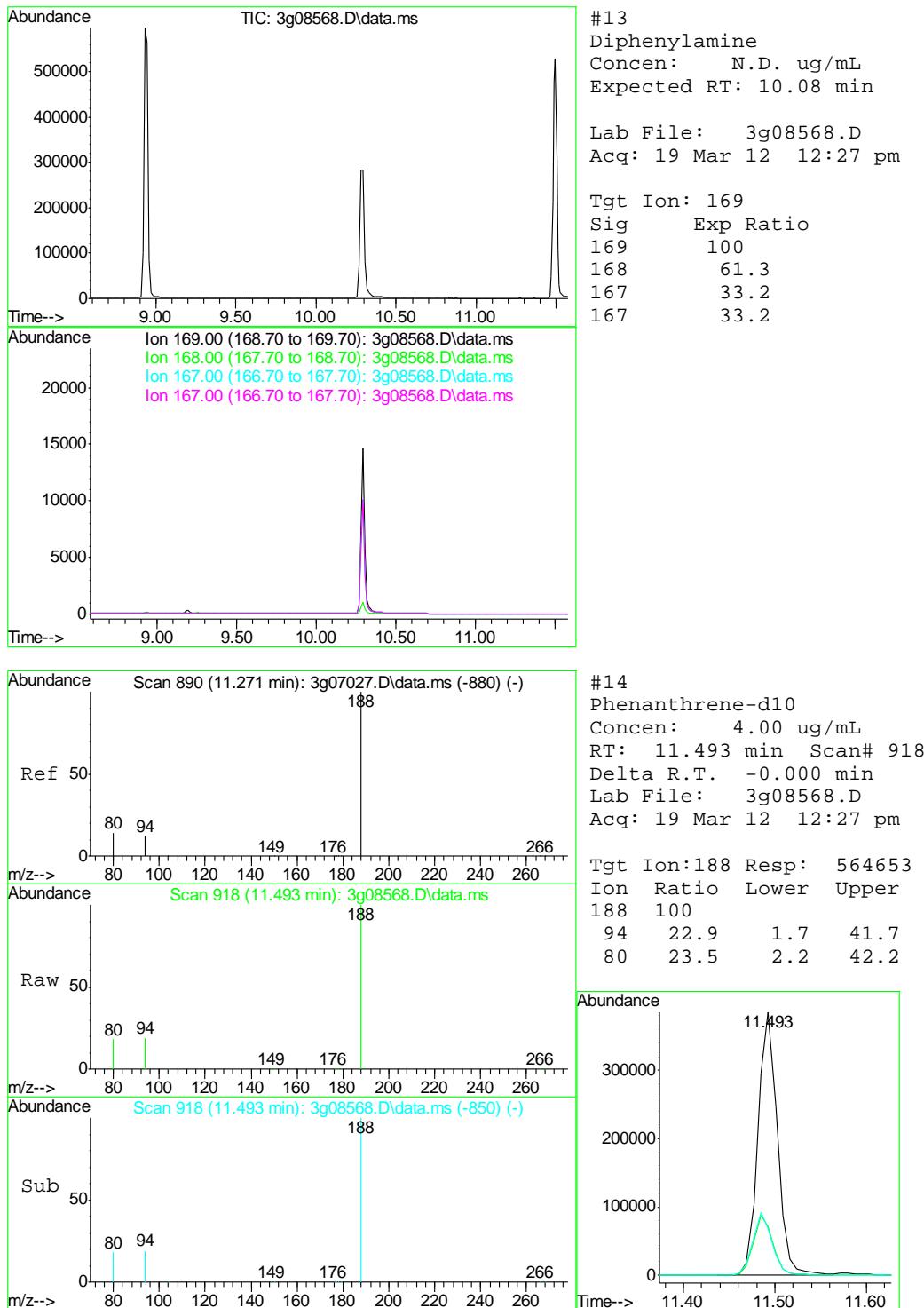


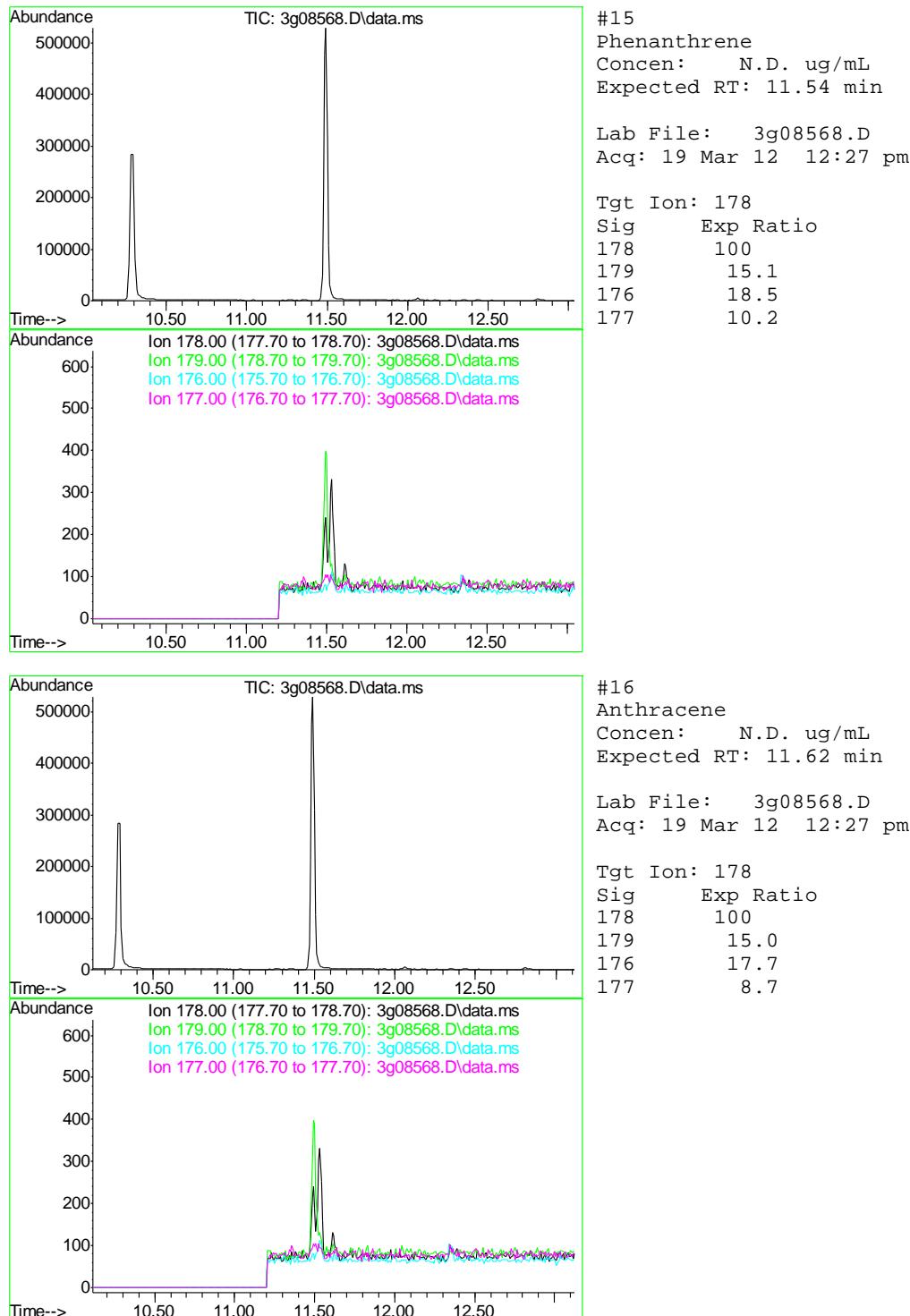


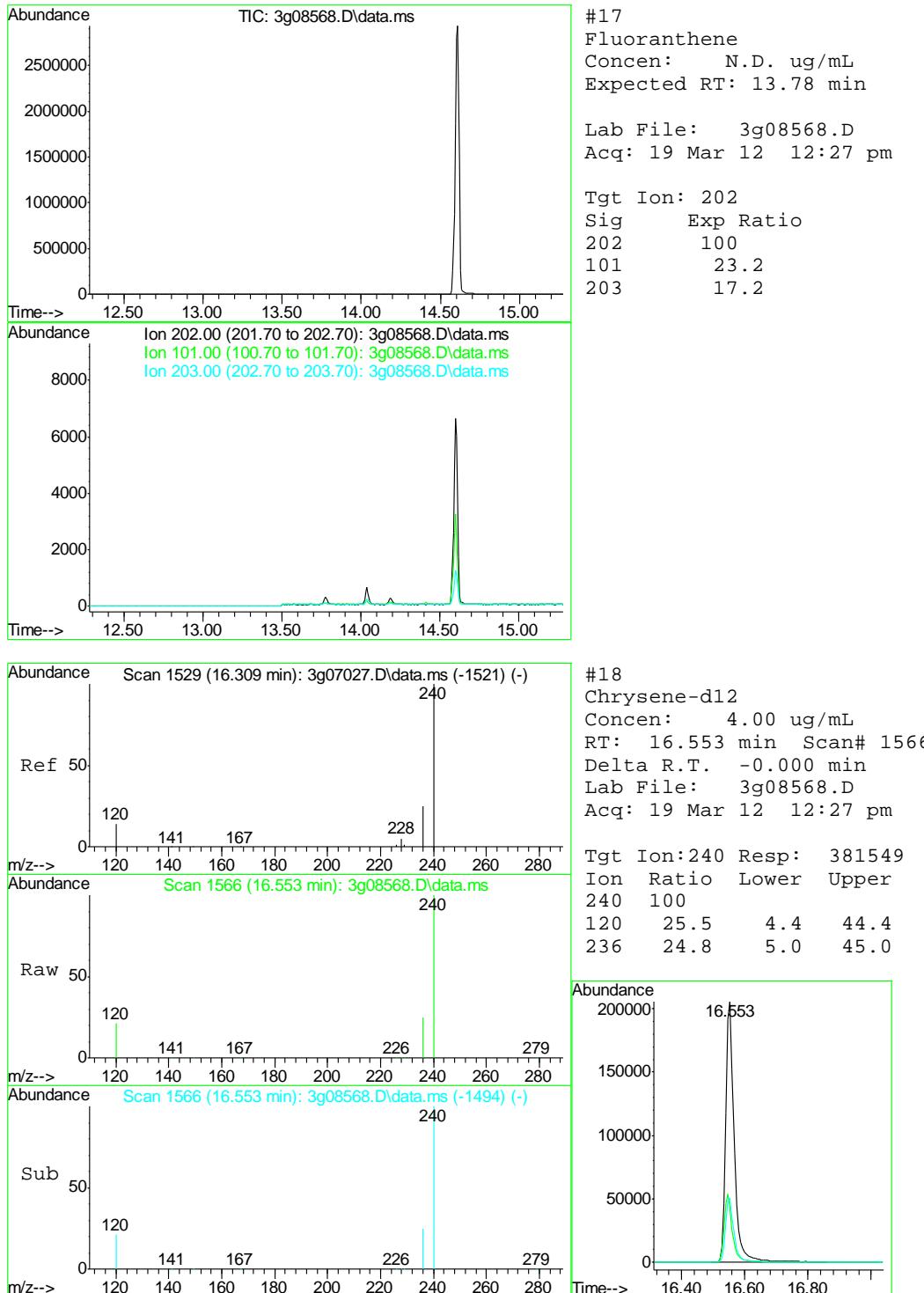


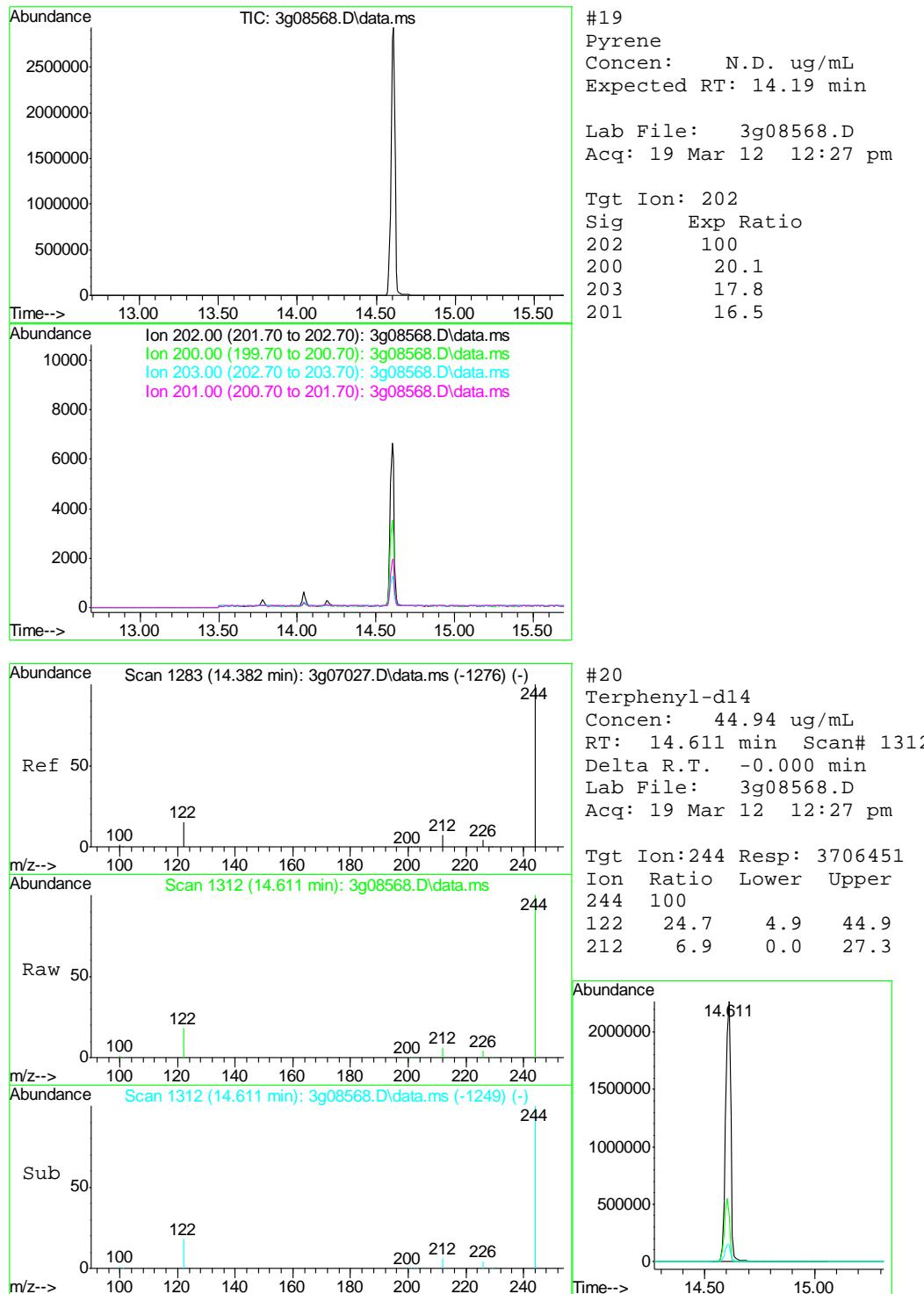


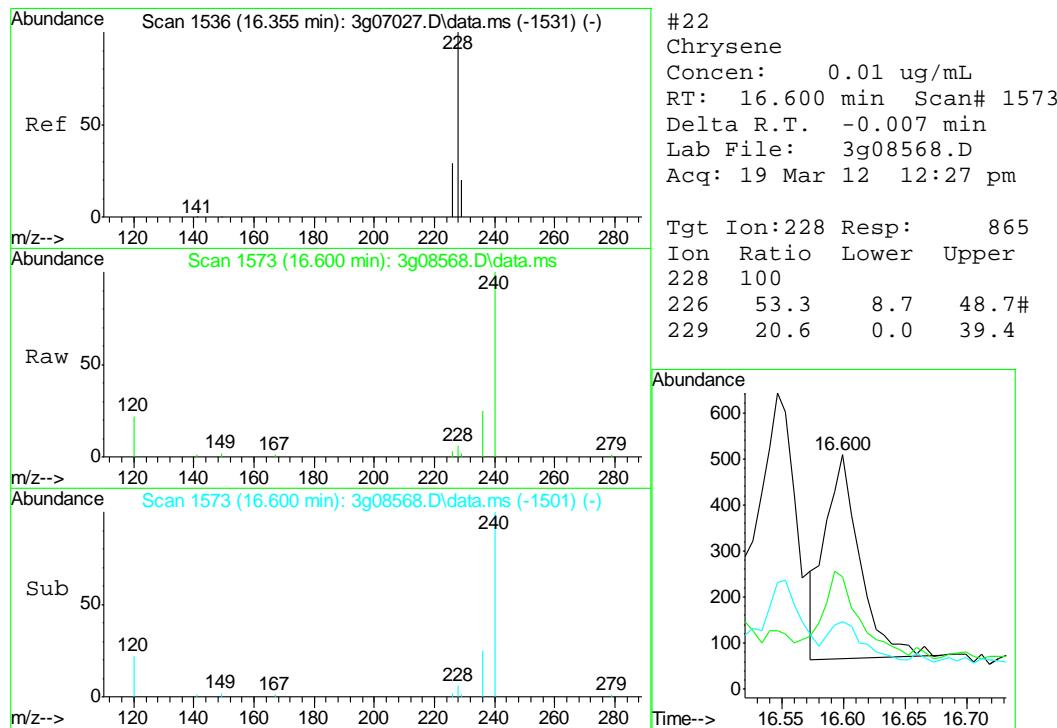
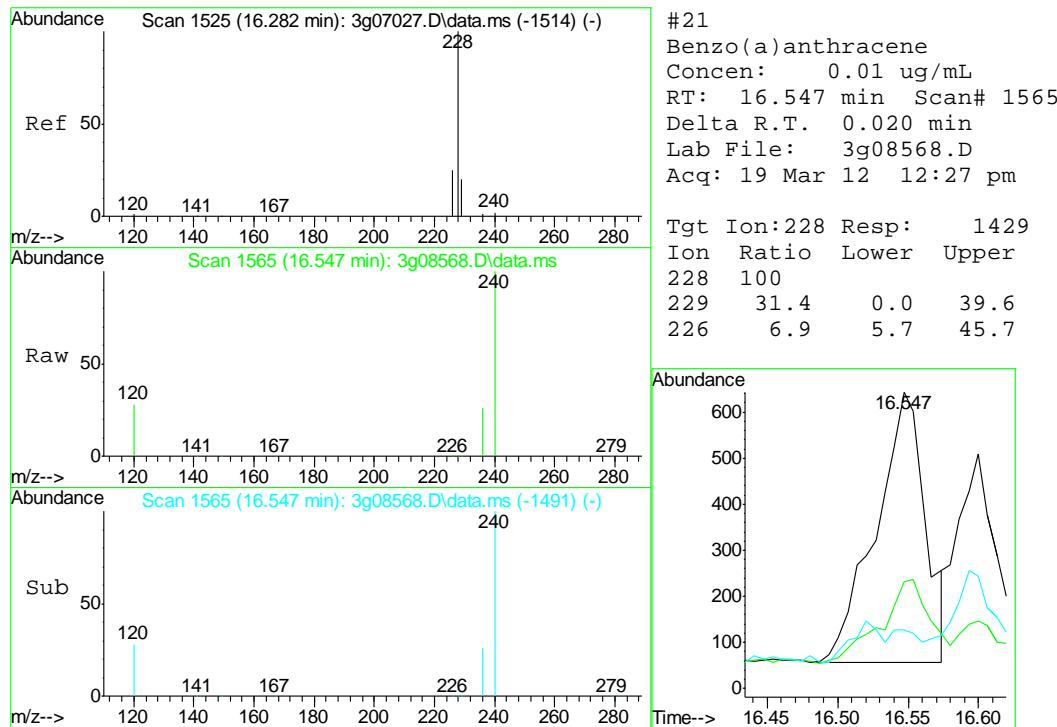


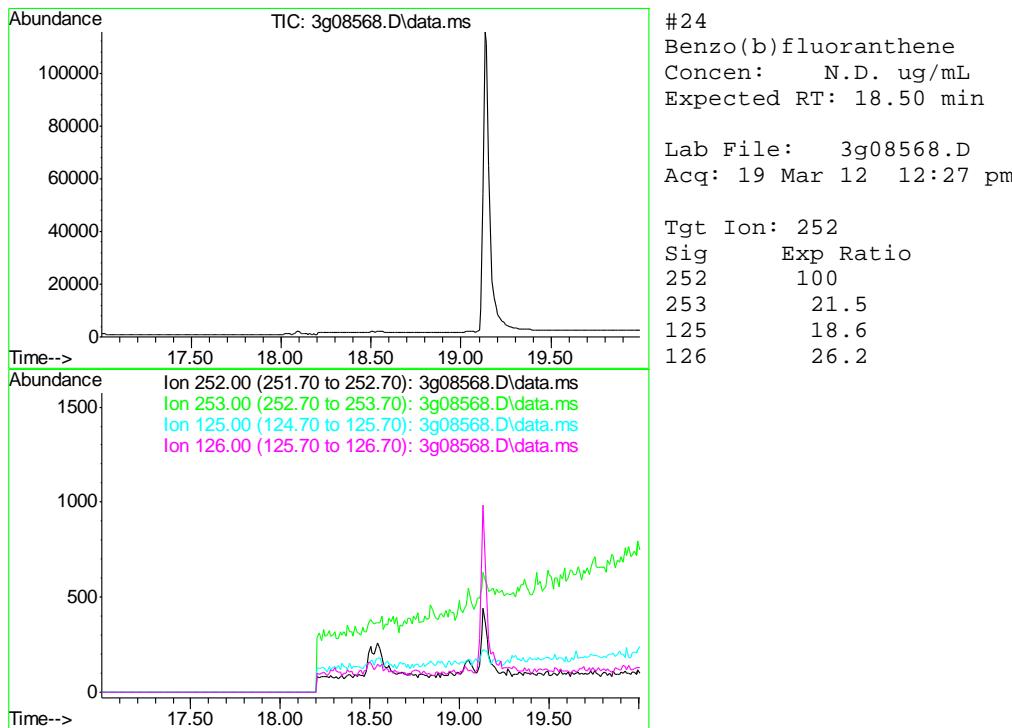
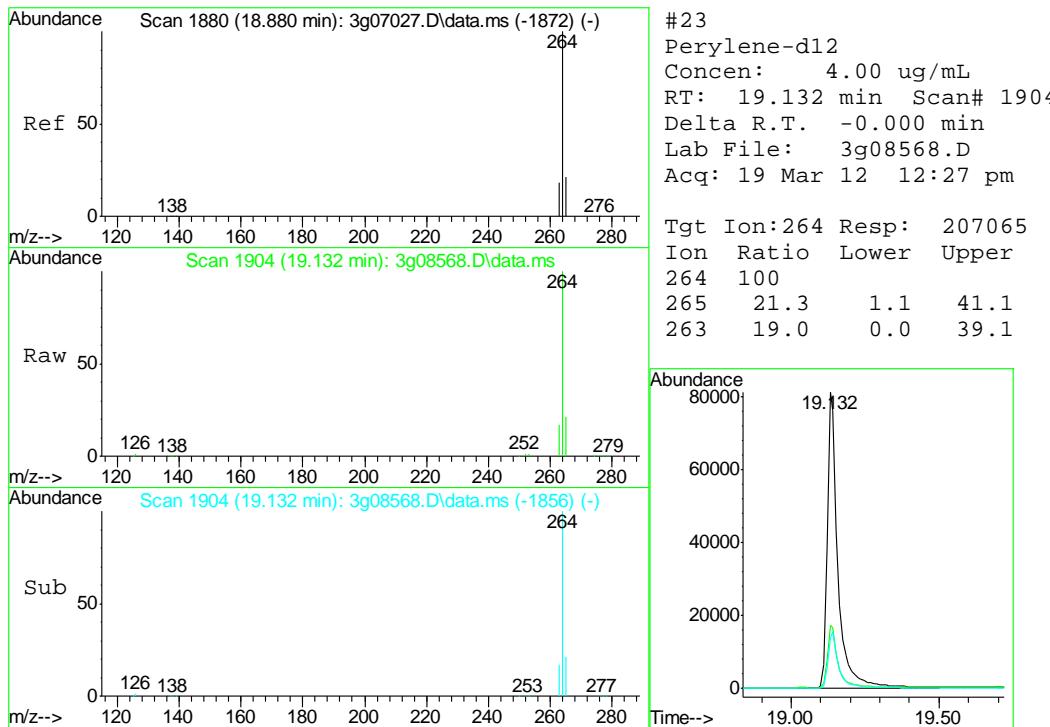


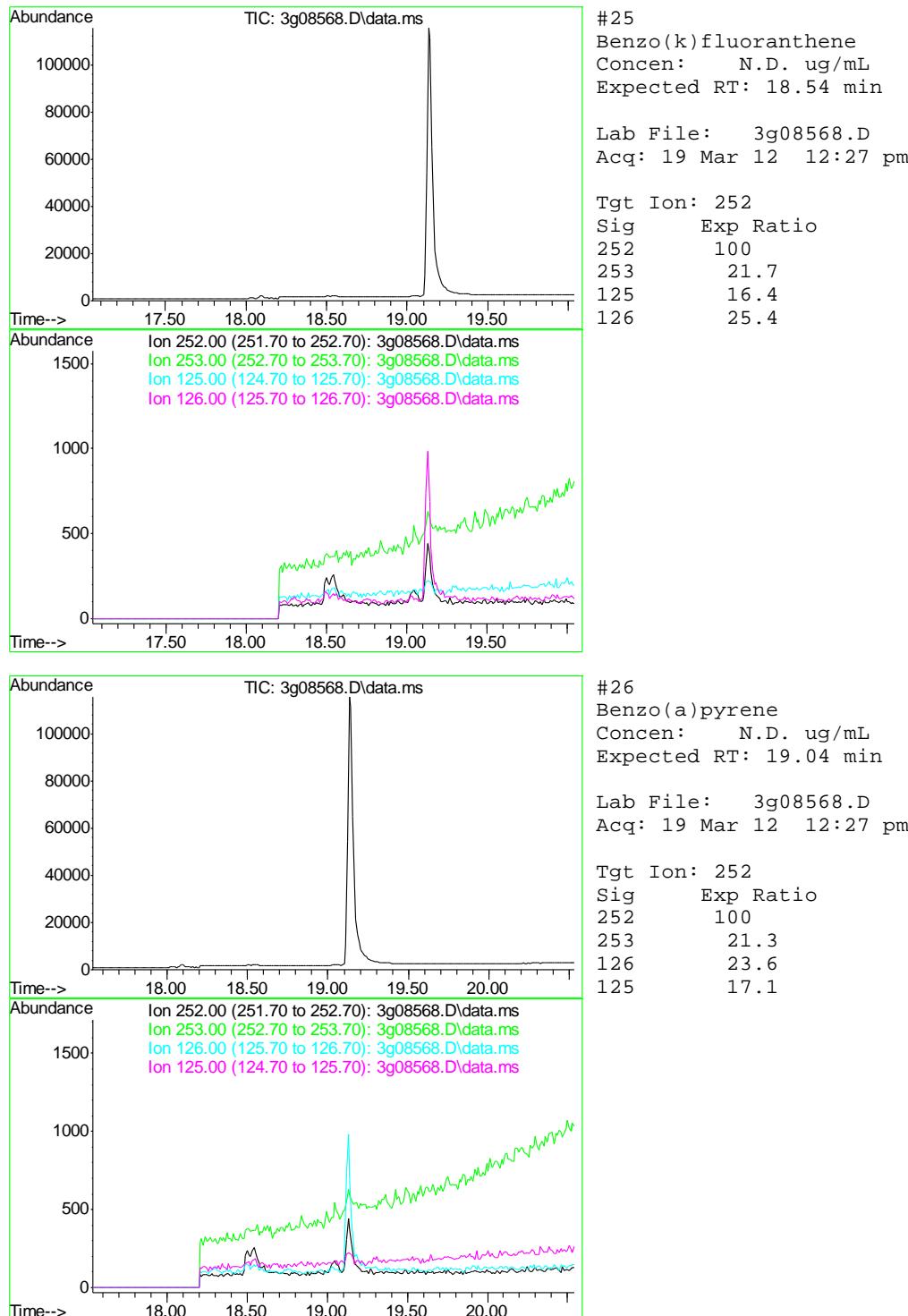


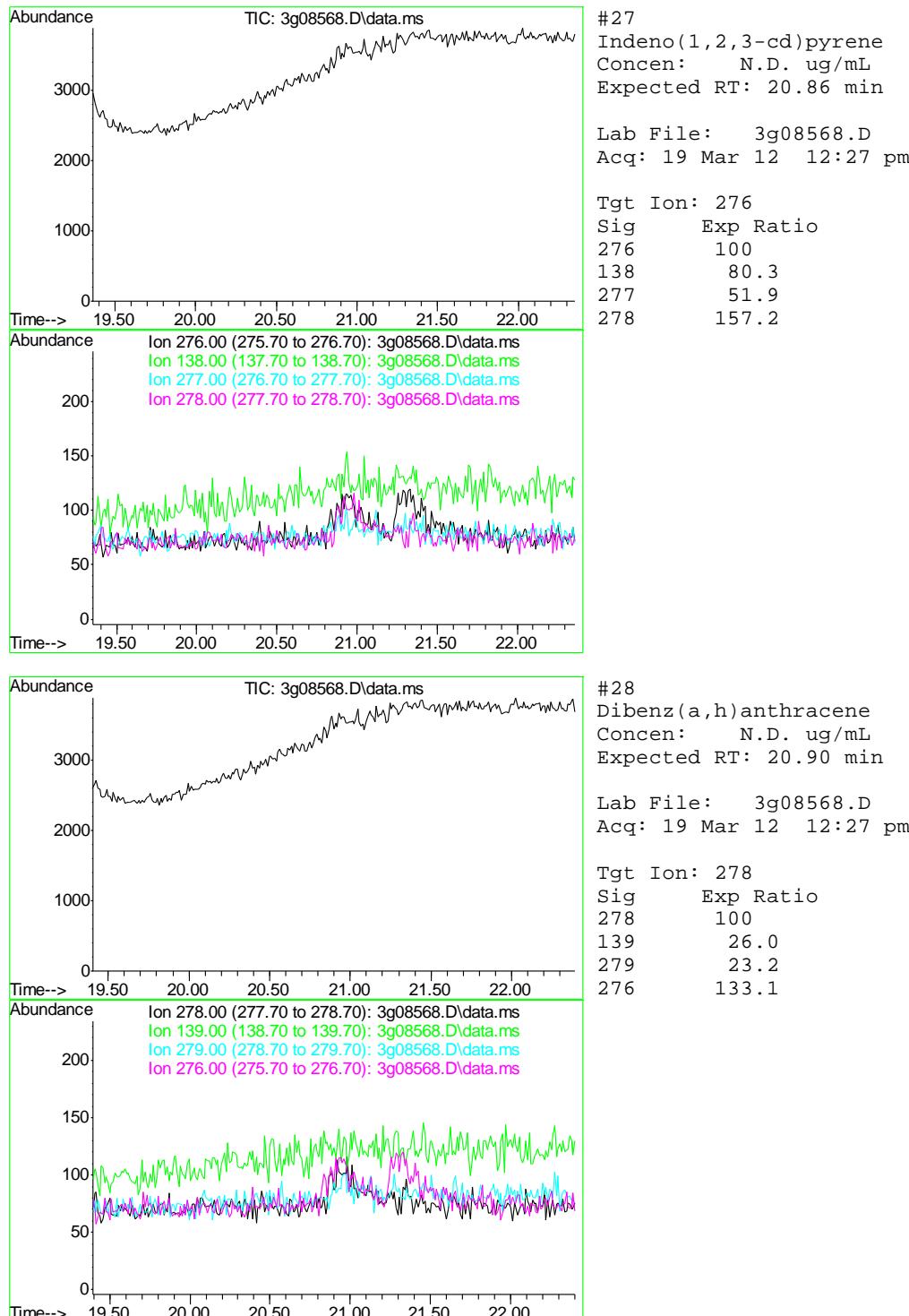


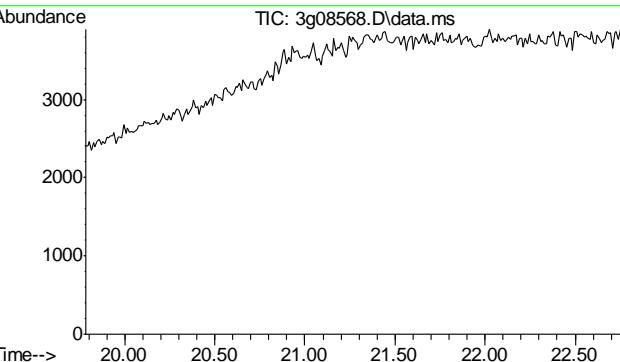








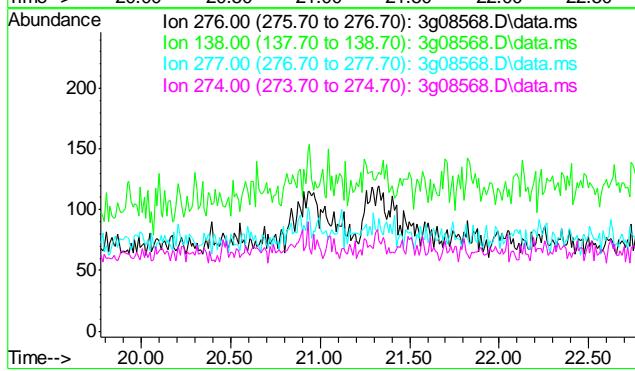




#29  
Benzo(g,h,i)perylene  
Concen: N.D. ug/mL  
Expected RT: 21.28 min

Lab File: 3g08568.D  
Acq: 19 Mar 12 12:27 pm

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





## 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:** D32747  
**Account:** XTOKWR XTO Energy  
**Project:** PCU T35X-2G

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
GGB859-MB	GB15306.D	1	03/15/12	SK	n/a	n/a	GGB859

The QC reported here applies to the following samples:

**Method:** SW846 8015B

D32747-1

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	104%      60-140%

9.1.1

9

## Blank Spike Summary

Page 1 of 1

Job Number: D32747

Account: XTOKWR XTO Energy

Project: PCU T35X-2G

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
GGB859-BS	GB15307.D	1	03/15/12	SK	n/a	n/a	GGB859

The QC reported here applies to the following samples:

Method: SW846 8015B

D32747-1

CAS No.	Compound	Spike mg/kg	BSP mg/kg	BSP %	Limits
	TPH-GRO (C6-C10)	110	114	104	70-130

CAS No.	Surrogate Recoveries	BSP	Limits
120-82-1	1,2,4-Trichlorobenzene	115%	60-140%

9.2.1

9

# Matrix Spike/Matrix Spike Duplicate Summary

Page 1 of 1

Job Number: D32747

Account: XTOKWR XTO Energy

Project: PCU T35X-2G

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
D32747-1MS	GB15309.D	1	03/15/12	SK	n/a	n/a	GGB859
D32747-1MSD	GB15310.D	1	03/15/12	SK	n/a	n/a	GGB859
D32747-1	GB15308.D	1	03/15/12	SK	n/a	n/a	GGB859

The QC reported here applies to the following samples:

Method: SW846 8015B

D32747-1

CAS No.	Compound	D32747-1		Spike	MS	MS	MSD	MSD	RPD	Limits Rec/RPD
		mg/kg	Q	mg/kg	mg/kg	%	mg/kg	%		
	TPH-GRO (C6-C10)	ND		124	126	101	128	103	2	70-130/30

CAS No.	Surrogate Recoveries	MS	MSD	D32747-1	Limits
120-82-1	1,2,4-Trichlorobenzene	109%	109%	96%	60-140%

9.3.1

9



## GC Volatiles

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Raw Data

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Judy Nelson  
 03/16/12 13:47

## Quantitation Report (QT Reviewed)

Signal #1 : Y:\1\DATA\031512\GB15308.D\FID1A.CH Vial: 6  
 Signal #2 : Y:\1\DATA\031512\GB15308.D\FID2B.CH  
 Acq On : 15 Mar 2012 9:26 pm Operator: StephK  
 Sample : D32747-1, 50X Inst : GC/MS Ins  
 Misc : GC2679,GGB859,5.007,,100,5,1 Multiplr: 1.00  
 IntFile Signal #1: TVH1.E IntFile Signal #2: FB2.E  
 Quant Time: Mar 16 07:16:01 2012 Quant Results File: TB851GB851SOIL.RES

Quant Method : C:\MSDCHEM\1...\TB851GB851SOIL.M (Chemstation Integrator)  
 Title : 8015B/8021B TVH/BTEX  
 Last Update : Fri Mar 16 07:15:36 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	14.35	2905995	96.330 %	m
10) S	1,2,4-Trichlorobenzene (P)	14.35	23623976	101.512 %	

Target Compounds

1) H	TVH-Gasoline	7.26	6301119	<MDL	mg/L
4) T	Methyl-t-butyl-ether	0.00	0	N.D.	ug/L d
5) T	Benzene	4.12	118474	0.214	ug/L
6) T	Toluene	7.63	517549	0.945	ug/L
7) T	Ethylbenzene	10.26	107345	0.235	ug/L
8) T	m,p-Xylene	10.44	648782	1.159	ug/L
9) T	o-Xylene	10.95	138041	0.301	ug/L
11) T	Naphthalene	14.53	1026254	3.928	ug/L

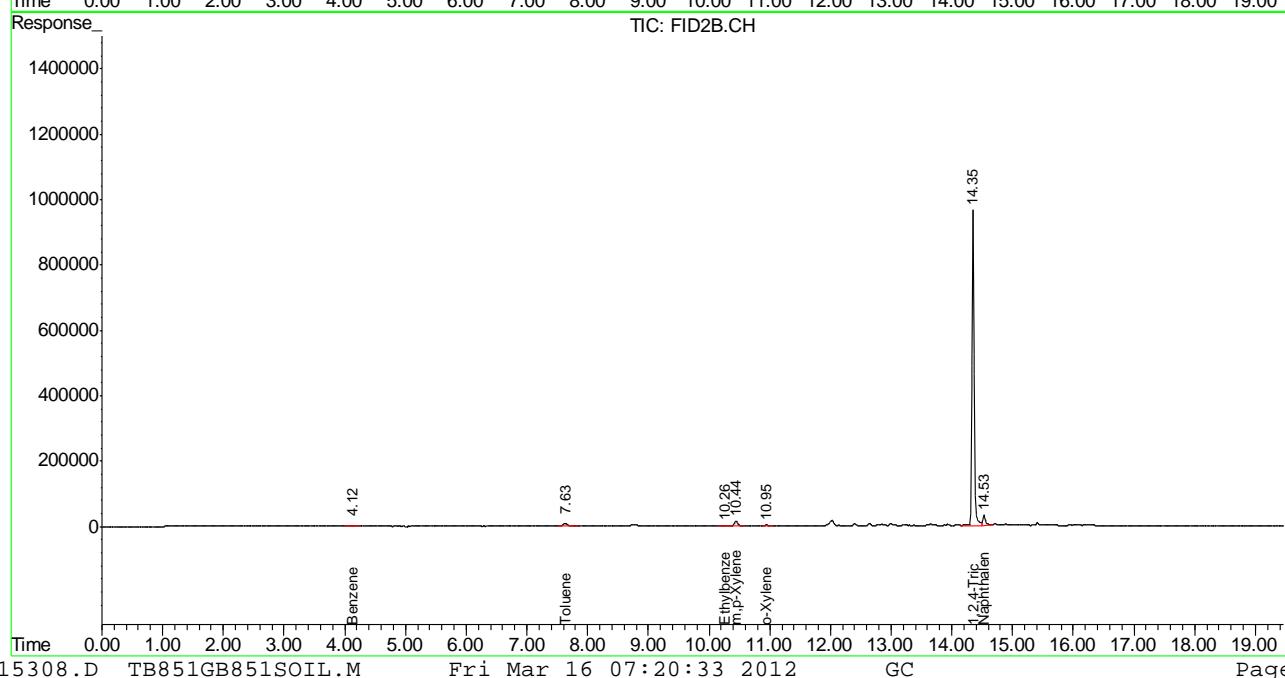
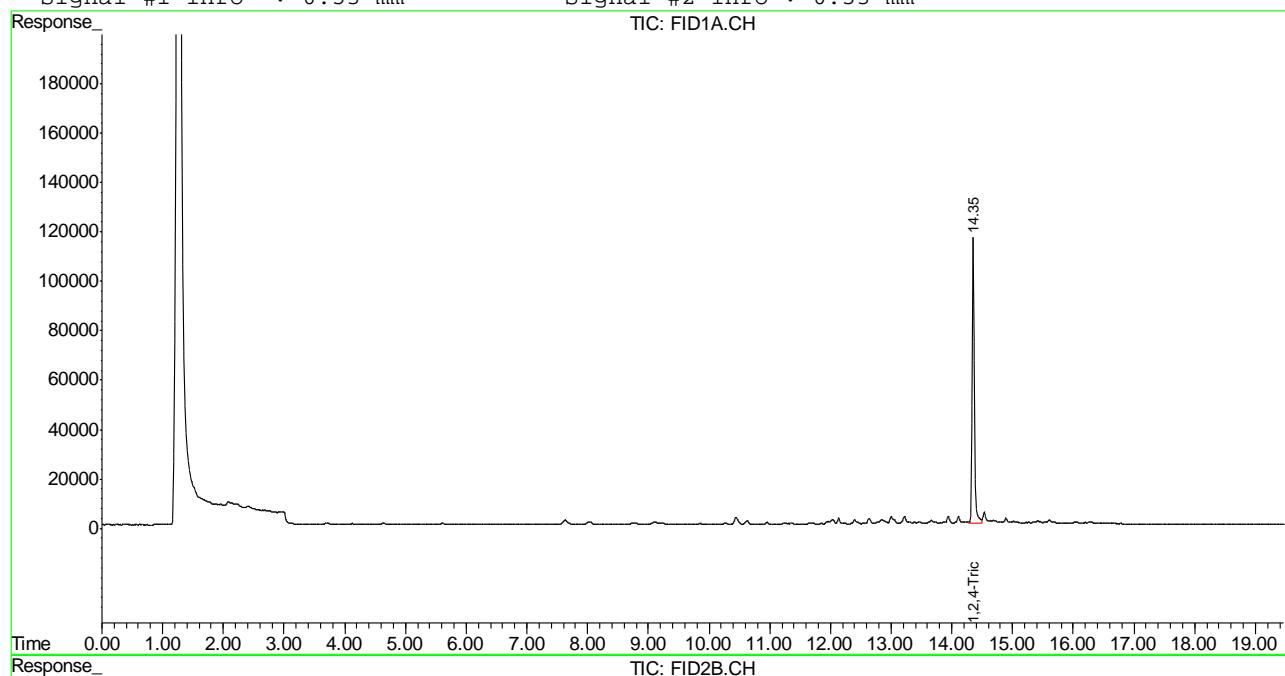
-----  
 (f)=RT Delta > 1/2 Window (m)=manual int.  
 GB15308.D TB851GB851SOIL.M Fri Mar 16 07:20:33 2012 GC

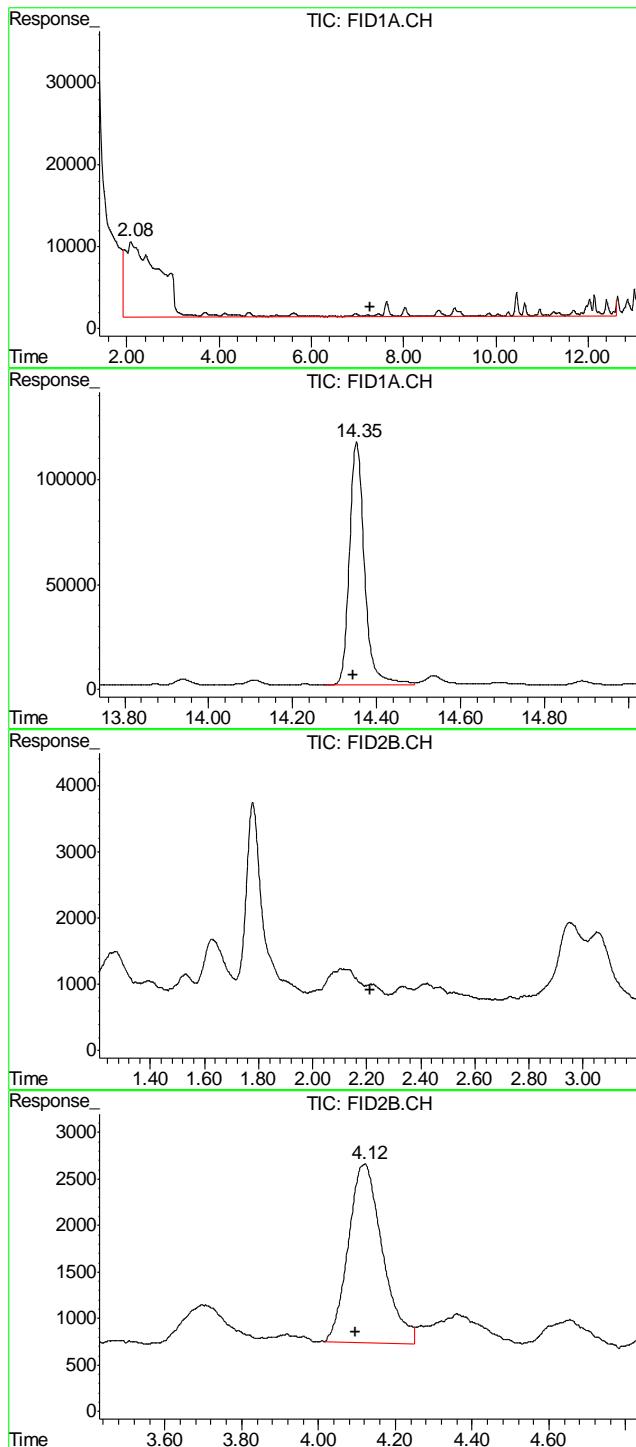
## Quantitation Report (QT Reviewed)

Signal #1 : Y:\1\DATA\031512\GB15308.D\FID1A.CH Vial: 6  
 Signal #2 : Y:\1\DATA\031512\GB15308.D\FID2B.CH  
 Acq On : 15 Mar 2012 9:26 pm Operator: StephK  
 Sample : D32747-1, 50X Inst : GC/MS Ins  
 Misc : GC2679,GGB859,5.007,,100,5,1 Multiplr: 1.00  
 IntFile Signal #1: TVH1.E IntFile Signal #2: FB2.E  
 Quant Time: Mar 16 7:19 2012 Quant Results File: TB851GB851SOIL.RES

Quant Method : C:\MSDCHEM\1...\TB851GB851SOIL.M (Chemstation Integrator)  
 Title : 8015B/8021B TVH/BTEX  
 Last Update : Fri Mar 16 07:15:36 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: 6301119  
 Conc: N.D.

## #2 1,2,4-Trichlorobenzene

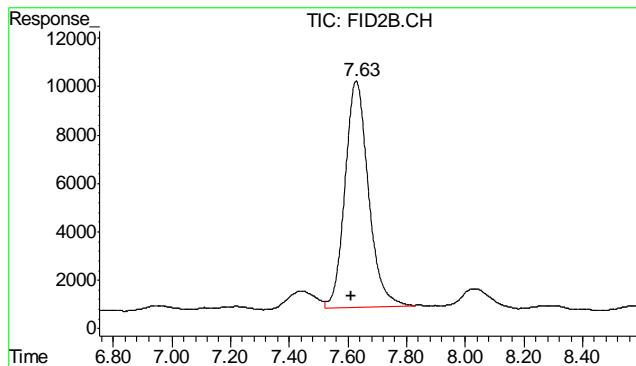
R.T.: 14.352 min  
 Delta R.T.: 0.007 min  
 Response: 2905995  
 Conc: 96.33 % m

## #4 Methyl-t-butyl-ether

R.T.: 0.000 min  
 Exp R.T.: 2.211 min  
 Response: 0  
 Conc: N.D.

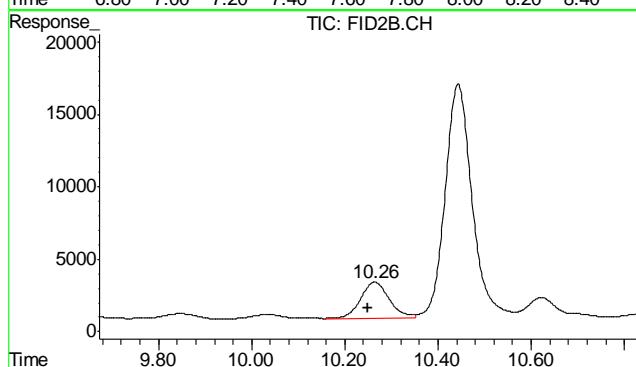
## #5 Benzene

R.T.: 4.122 min  
 Delta R.T.: 0.026 min  
 Response: 118474  
 Conc: 0.21 ug/L



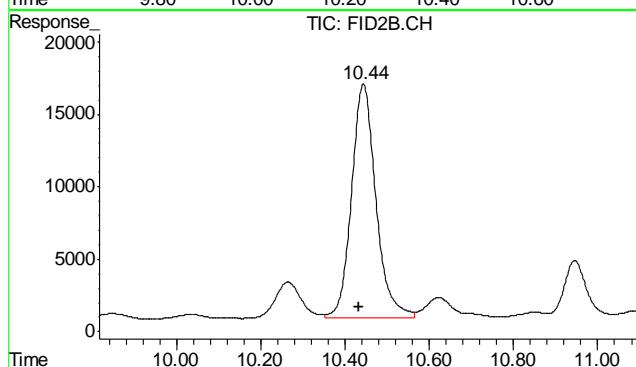
#6 Toluene

R.T.: 7.628 min  
 Delta R.T.: 0.016 min  
 Response: 517549  
 Conc: 0.95 ug/L



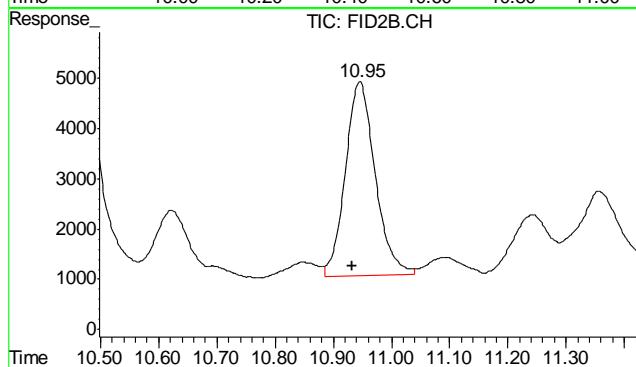
#7 Ethylbenzene

R.T.: 10.264 min  
 Delta R.T.: 0.013 min  
 Response: 107345  
 Conc: 0.23 ug/L



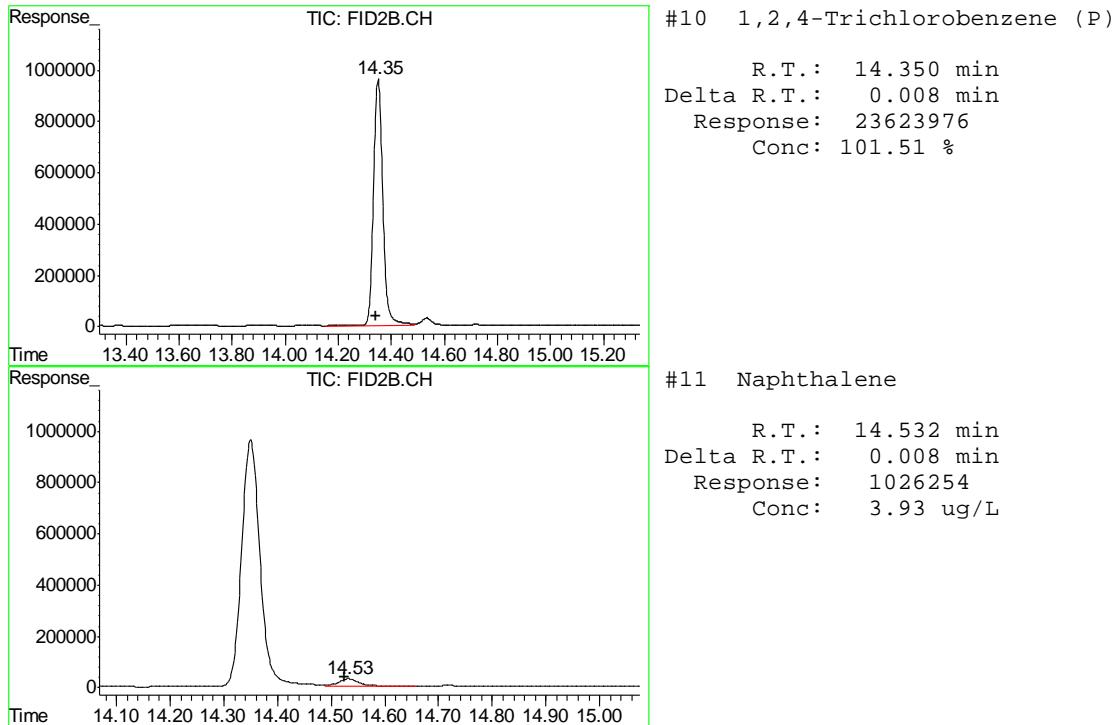
#8 m,p-Xylene

R.T.: 10.443 min  
 Delta R.T.: 0.010 min  
 Response: 648782  
 Conc: 1.16 ug/L



#9 o-Xylene

R.T.: 10.946 min  
 Delta R.T.: 0.013 min  
 Response: 138041  
 Conc: 0.30 ug/L



Judy Nelson  
 03/16/12 13:47

## Quantitation Report (QT Reviewed)

Signal #1 : Y:\1\DATA\031512\GB15306.D\FID1A.CH Vial: 4  
 Signal #2 : Y:\1\DATA\031512\GB15306.D\FID2B.CH  
 Acq On : 15 Mar 2012 8:15 pm Operator: StephK  
 Sample : MB Inst : GC/MS Ins  
 Misc : GC2679,GGB859,5.000,,100,5,1 Multiplr: 1.00  
 IntFile Signal #1: TVH1.E IntFile Signal #2: FB2.E  
 Quant Time: Mar 16 07:15:53 2012 Quant Results File: TB851GB851SOIL.RES

Quant Method : C:\MSDCHEM\1...\TB851GB851SOIL.M (Chemstation Integrator)  
 Title : 8015B/8021B TVH/BTEX  
 Last Update : Fri Mar 16 07:15:36 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	3134910	103.918 %	m
10) S	1,2,4-Trichlorobenzene (P)	14.33	25016008	107.494 %	

**Target Compounds**

1) H	TVH-Gasoline	7.26	5799283	<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	7.61	268803	0.491	ug/L
7) T	Ethylbenzene	0.00	0	N.D.	ug/L d
8) T	m,p-Xylene	10.43	247992	0.443	ug/L
9) T	o-Xylene	10.93	88682	0.193	ug/L
11) T	Naphthalene	14.51	335362	1.283	ug/L

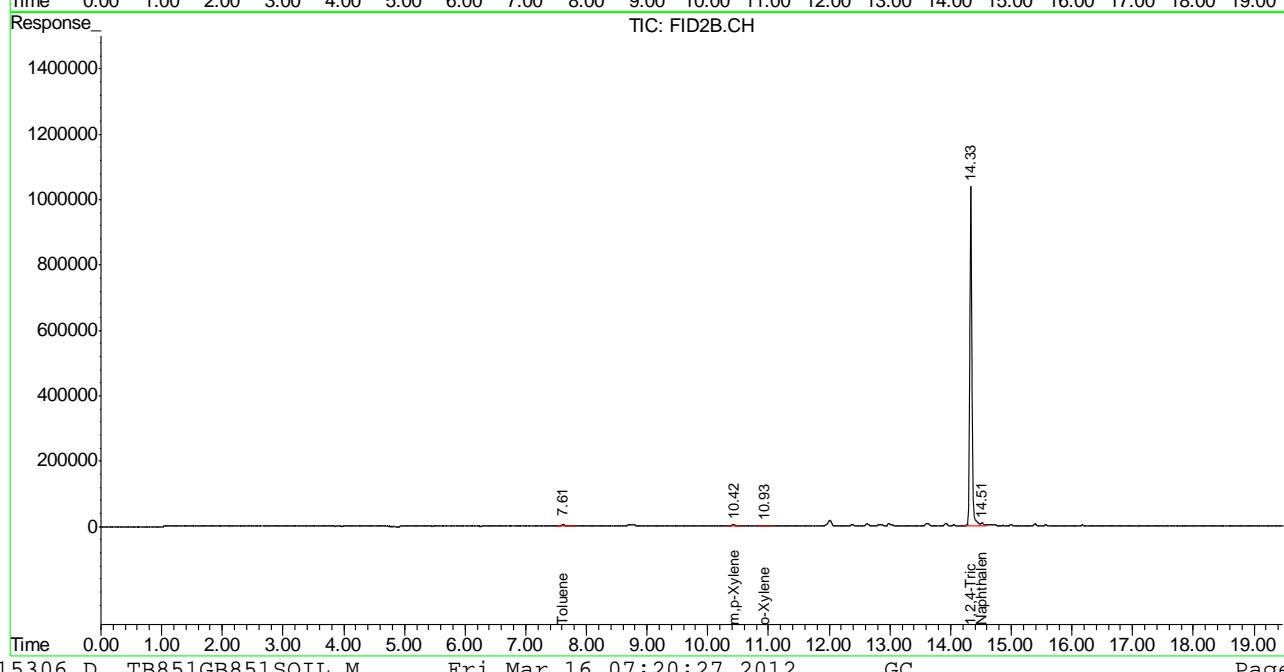
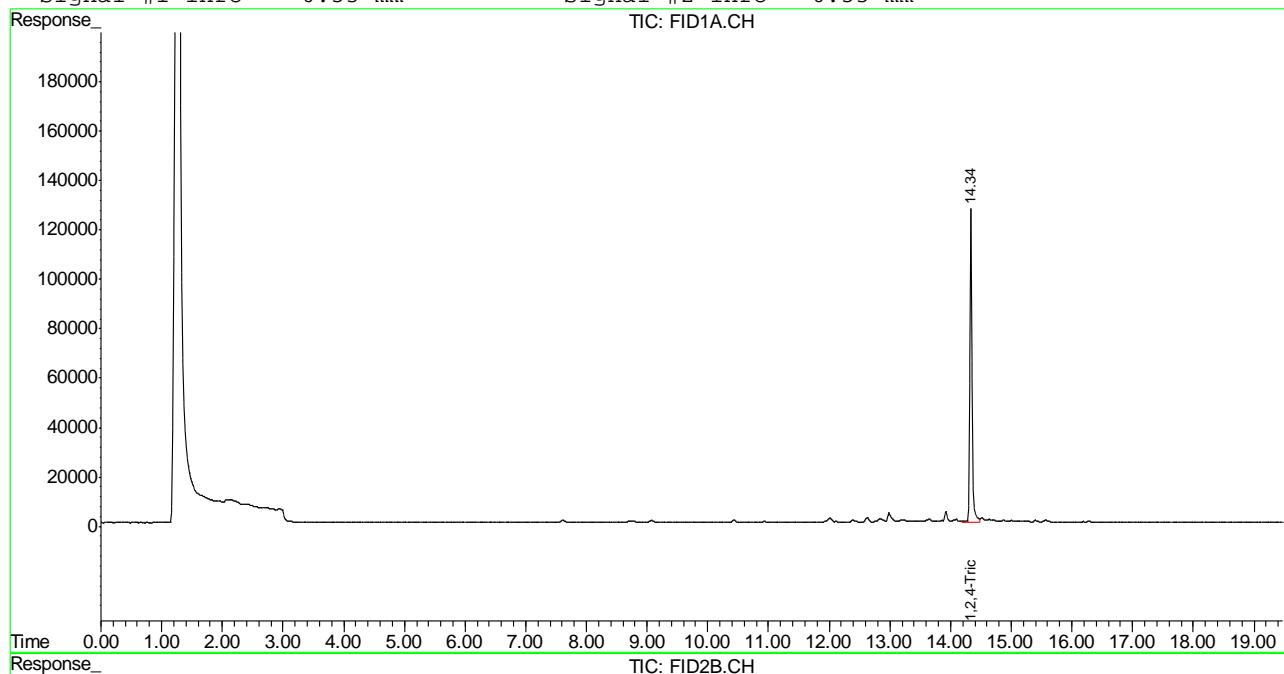
(f)=RT Delta > 1/2 Window (m)=manual int.  
 GB15306.D TB851GB851SOIL.M Fri Mar 16 07:20:27 2012 GC

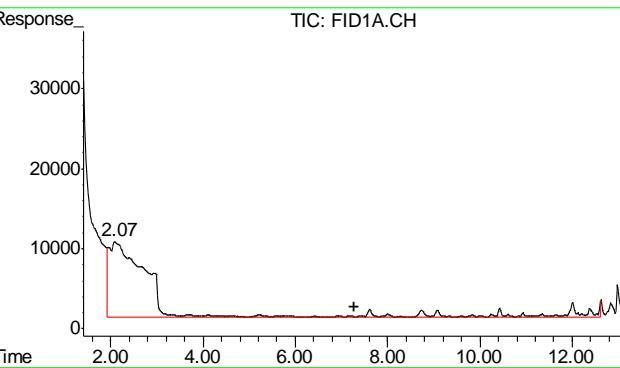
## Quantitation Report (QT Reviewed)

Signal #1 : Y:\1\DATA\031512\GB15306.D\FID1A.CH Vial: 4  
 Signal #2 : Y:\1\DATA\031512\GB15306.D\FID2B.CH  
 Acq On : 15 Mar 2012 8:15 pm Operator: StephK  
 Sample : MB Inst : GC/MS Ins  
 Misc : GC2679, GGB859, 5.000,,100,5,1 Multiplr: 1.00  
 IntFile Signal #1: TVH1.E IntFile Signal #2: FB2.E  
 Quant Time: Mar 16 7:18 2012 Quant Results File: TB851GB851SOIL.RES

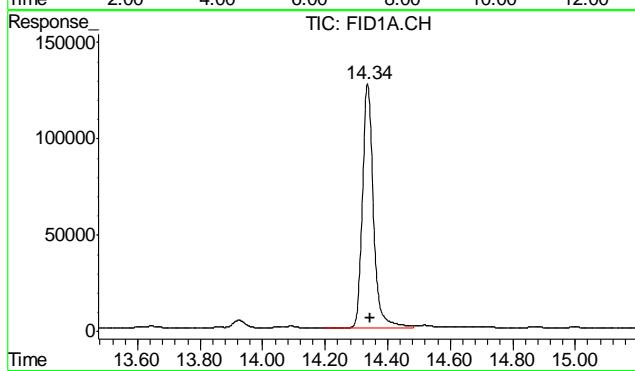
Quant Method : C:\MSDCHEM\1...\TB851GB851SOIL.M (Chemstation Integrator)  
 Title : 8015B/8021B TVH/BTEX  
 Last Update : Fri Mar 16 07:15:36 2012  
 Response via : Multiple Level Calibration  
 DataAcq Meth : TVB4.M

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

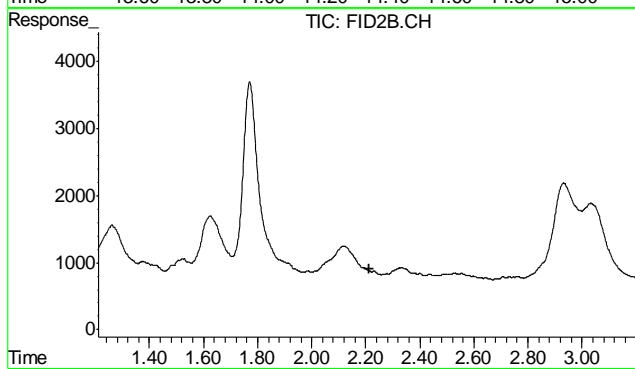




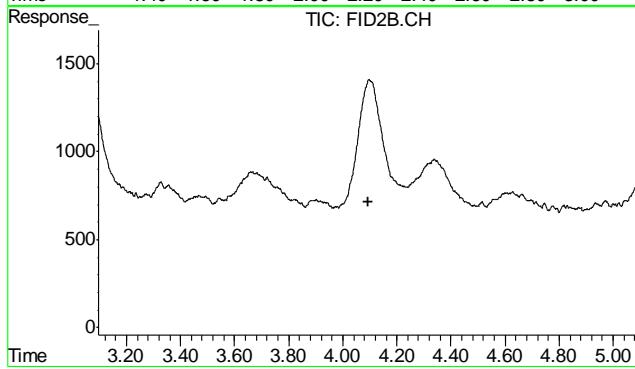
#1 TVH-Gasoline  
R.T.: 7.265 min  
Delta R.T.: 0.000 min  
Response: 5799283  
Conc: N.D.



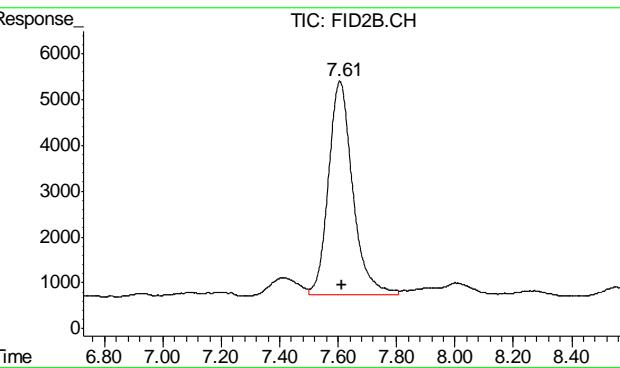
#2 1,2,4-Trichlorobenzene  
R.T.: 14.335 min  
Delta R.T.: -0.009 min  
Response: 3134910  
Conc: 103.92 % m



#4 Methyl-t-butyl-ether  
R.T.: 0.000 min  
Exp R.T. : 2.211 min  
Response: 0  
Conc: N.D.

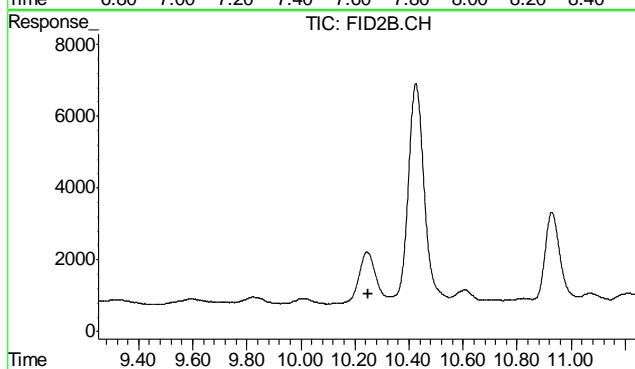


#5 Benzene  
R.T.: 0.000 min  
Exp R.T. : 4.096 min  
Response: 0  
Conc: N.D.



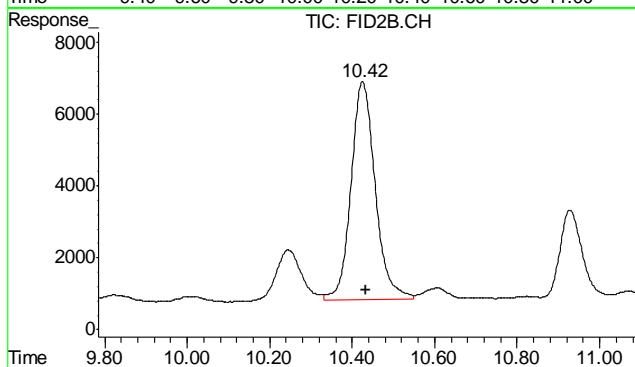
#6 Toluene

R.T.: 7.606 min  
Delta R.T.: -0.006 min  
Response: 268803  
Conc: 0.49 ug/L



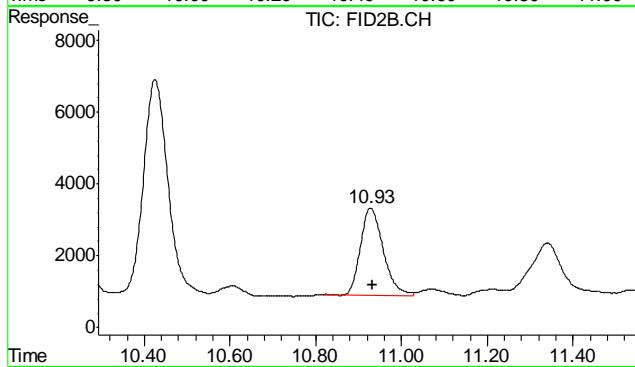
#7 Ethylbenzene

R.T.: 0.000 min  
Exp R.T. : 10.251 min  
Response: 0  
Conc: N.D.



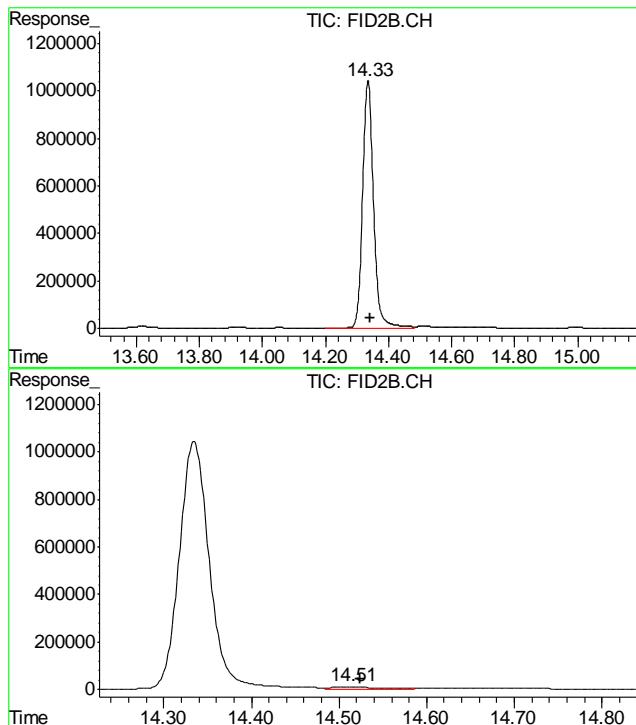
#8 m,p-Xylene

R.T.: 10.425 min  
Delta R.T.: -0.008 min  
Response: 247992  
Conc: 0.44 ug/L



#9 o-Xylene

R.T.: 10.928 min  
Delta R.T.: -0.005 min  
Response: 88682  
Conc: 0.19 ug/L



#10 1,2,4-Trichlorobenzene (P)

R.T.: 14.335 min  
Delta R.T.: -0.007 min  
Response: 25016008  
Conc: 107.49 %

#11 Naphthalene

R.T.: 14.515 min  
Delta R.T.: -0.009 min  
Response: 335362  
Conc: 1.28 ug/L



## 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: D32747

Account: XTOKWR XTO Energy

Project: PCU T35X-2G

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
OP5560-MB	FH002331.D	1	03/20/12	TR	03/19/12	OP5560	GFH123

The QC reported here applies to the following samples:

**Method:** SW846-8015B

D32747-1

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	86% 43-136%

11.11

11

## Blank Spike Summary

Page 1 of 1

Job Number: D32747

Account: XTOKWR XTO Energy

Project: PCU T35X-2G

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
OP5560-BS	FH002333.D	1	03/20/12	TR	03/19/12	OP5560	GFH123

The QC reported here applies to the following samples:

Method: SW846-8015B

D32747-1

CAS No.	Compound	Spike mg/kg	BSP mg/kg	BSP %	Limits
	TPH-DRO (C10-C28)	667	520	78	58-130

CAS No.	Surrogate Recoveries	BSP	Limits
84-15-1	o-Terphenyl	85%	43-136%

11.2.1  
11

# Matrix Spike/Matrix Spike Duplicate Summary

Page 1 of 1

Job Number: D32747

Account: XTOKWR XTO Energy

Project: PCU T35X-2G

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
OP5560-MS	FH002335.D	1	03/20/12	TR	03/19/12	OP5560	GFH123
OP5560-MSD	FH002337.D	1	03/20/12	TR	03/19/12	OP5560	GFH123
D32747-1	FH002339.D	1	03/20/12	TR	03/19/12	OP5560	GFH123

The QC reported here applies to the following samples:

Method: SW846-8015B

D32747-1

CAS No.	Compound	D32747-1		Spike	MS	MS	MSD	MSD	RPD	Limits Rec/RPD
		mg/kg	Q	mg/kg	mg/kg	%	mg/kg	%		
	TPH-DRO (C10-C28)	31.1		710	518	69	483	64	7	20-183/43
CAS No.	Surrogate Recoveries	MS	MSD	D32747-1		Limits				
84-15-1	o-Terphenyl	68%	68%	64%		43-136%				

11.3.1  
11



GC Semi-volatiles

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Raw Data

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12

## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\FH032012.SEC\  
 Data File : FH002339.D  
 Signal(s) : FID2B.ch  
 Acq On : 20 Mar 2012 6:00 pm  
 Operator : teder  
 Sample : D32747-1  
 Misc : OP5560,GFH123,30.12,,,2,1  
 ALS Vial : 57 Sample Multiplier: 1

Integration File: events.e  
 Quant Time: Mar 21 08:17:25 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	12.344	1098692658	641.625	ug/ml
<hr/>				
Target Compounds				
1) H TPH-DRO (C10-C28)	9.832	678039155	439.362	ug/ml
<hr/>				

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

(m)=manual int.

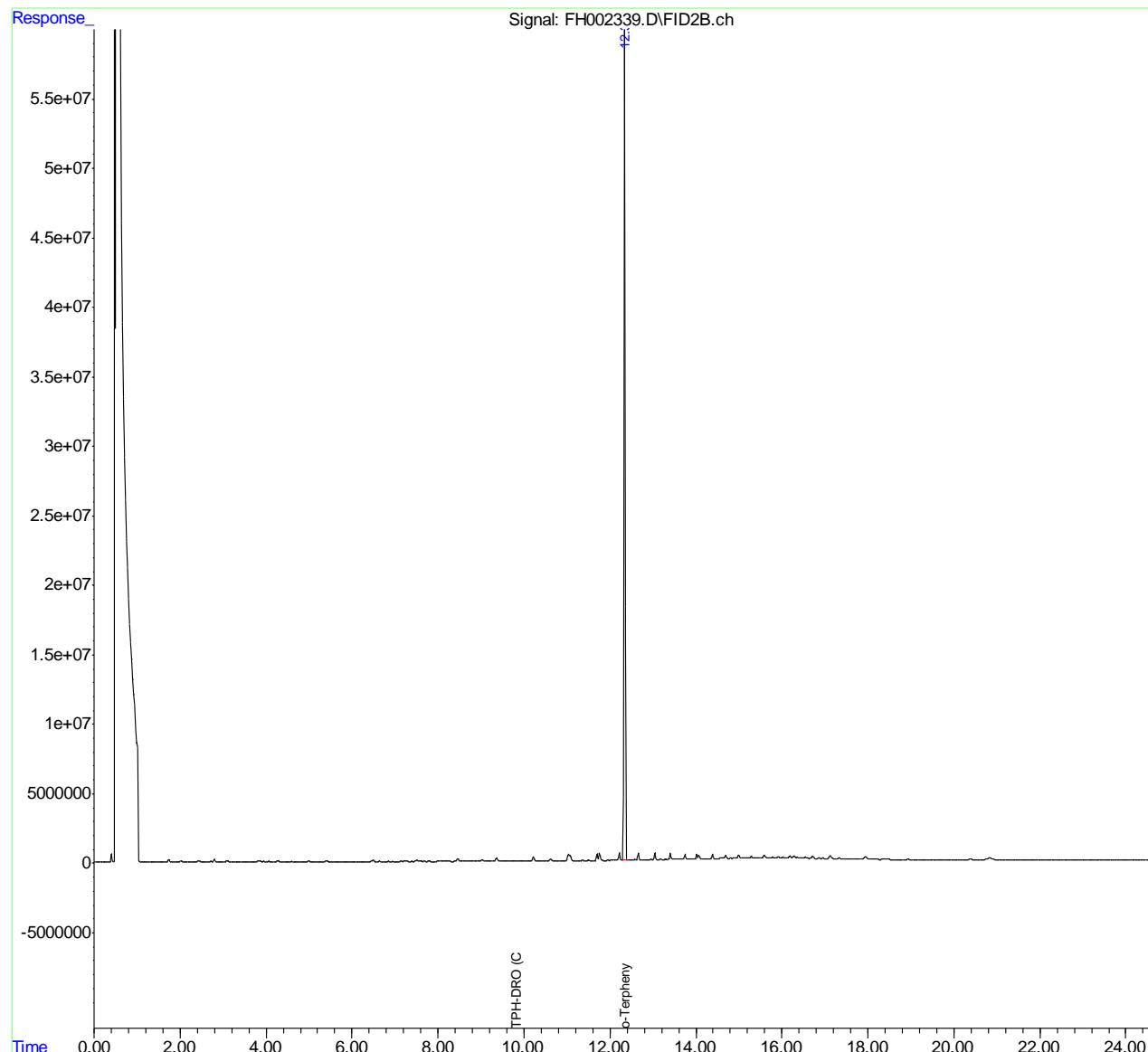
12.1.1  
12

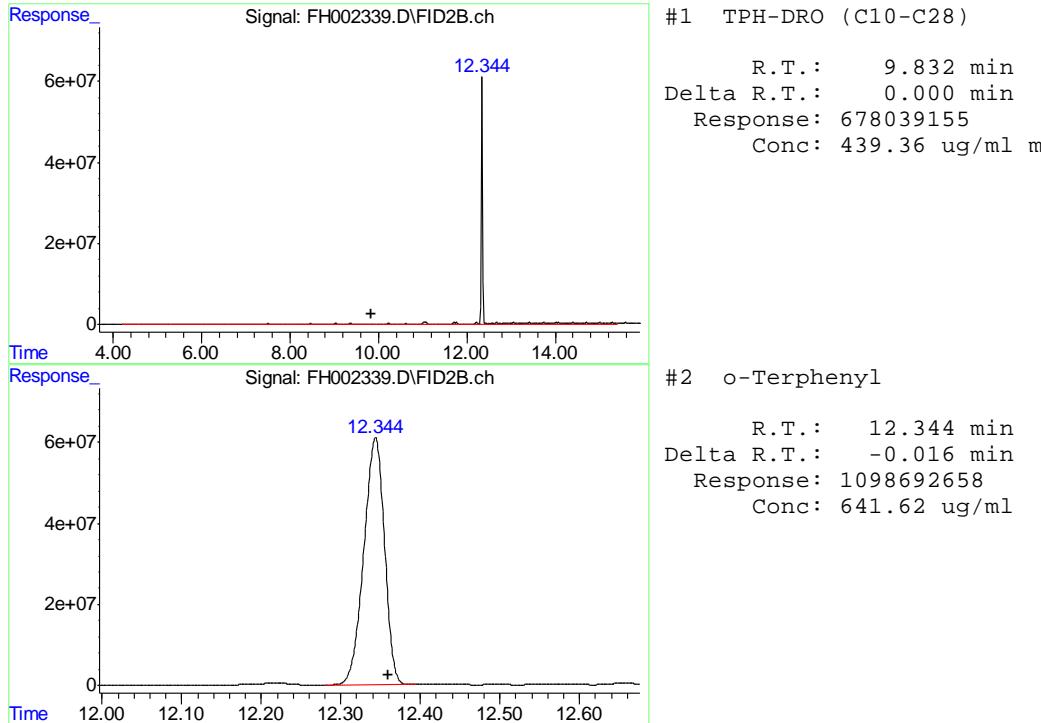
## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\FH032012.SEC\  
 Data File : FH002339.D  
 Signal(s) : FID2B.ch  
 Acq On : 20 Mar 2012 6:00 pm  
 Operator : tedr  
 Sample : D32747-1  
 Misc : OP5560,GFH123,30.12,,,2,1  
 ALS Vial : 57 Sample Multiplier: 1

Integration File: events.e  
 Quant Time: Mar 21 08:17:25 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\FH032012.SEC\  
 Data File : FH002331.D  
 Signal(s) : FID2B.ch  
 Acq On : 20 Mar 2012 3:38 pm  
 Operator : teder  
 Sample : OP5560-MB  
 Misc : OP5560,GFH123,30.00,,,2,1  
 ALS Vial : 53 Sample Multiplier: 1

Integration File: events.e  
 Quant Time: Mar 20 16:04:08 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	12.346	1475787544	861.844	ug/ml
<hr/>				
Target Compounds				
1) H TPH-DRO (C10-C28)	9.832	43097947	27.927	ug/ml
<hr/>				

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

12.2.1

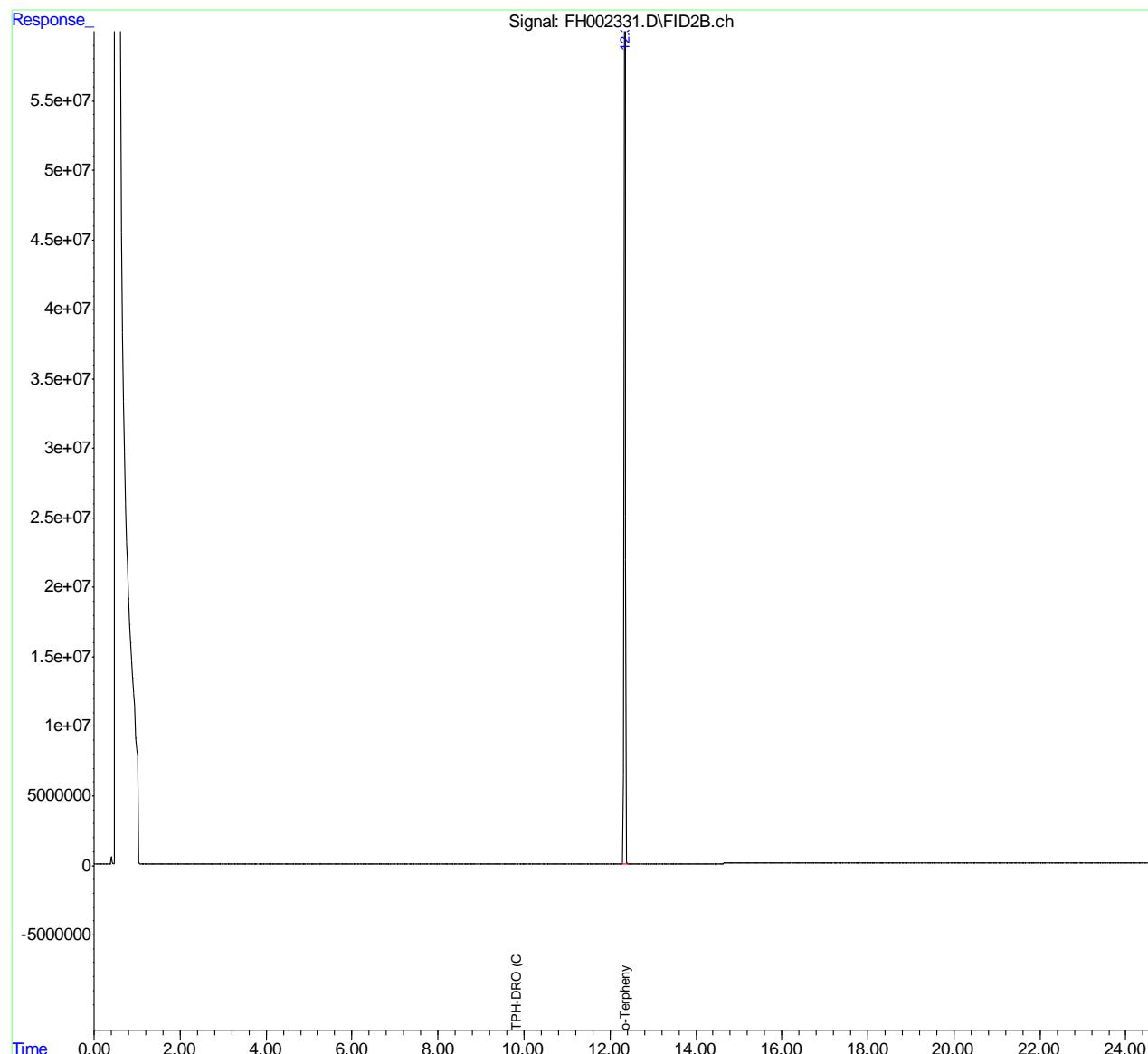
12

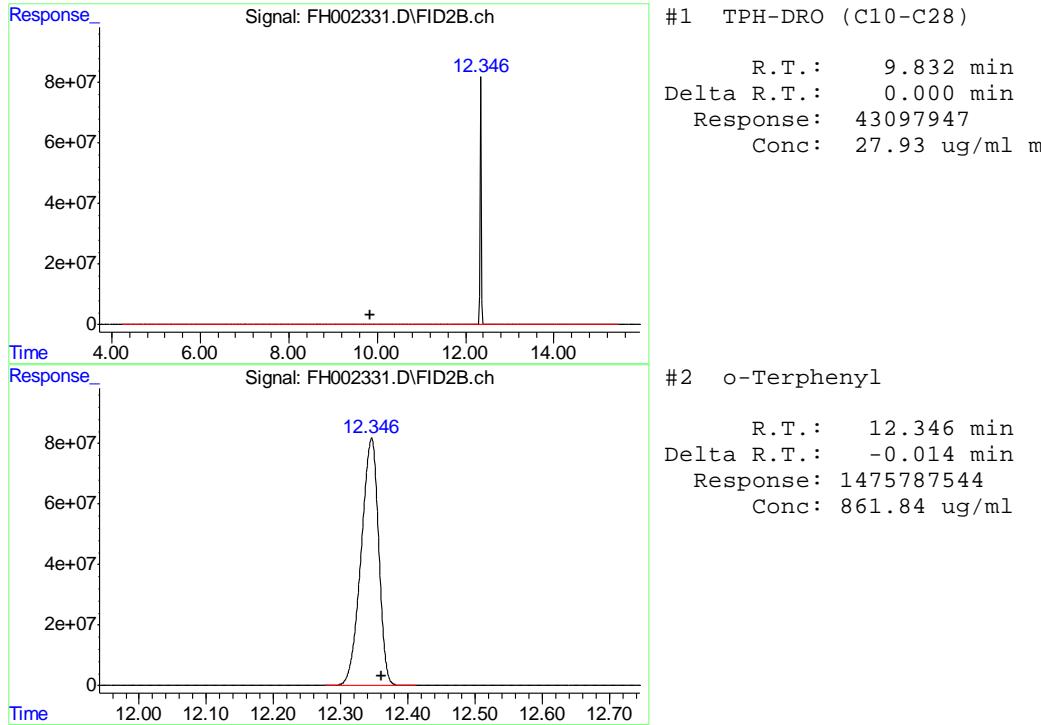
## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\FH032012.SEC\  
 Data File : FH002331.D  
 Signal(s) : FID2B.ch  
 Acq On : 20 Mar 2012 3:38 pm  
 Operator : tedr  
 Sample : OP5560-MB  
 Misc : OP5560,GFH123,30.00,,,2,1  
 ALS Vial : 53 Sample Multiplier: 1

Integration File: events.e  
 Quant Time: Mar 20 16:04:08 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

---

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: D32747  
Account: XTOKRWR - XTO Energy  
Project: PCU T35X-2G

QC Batch ID: MP7078  
Matrix Type: SOLID

Methods: SW846 6010C  
Units: mg/kg

Prep Date:

03/16/12

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	0.060	<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	0.19	<1.0
Iron	7.0	.34	2		
Lead	5.0	.16	.21	0.20	<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	0.040	<3.0
Phosphorus	10	1.1	1.2		
Potassium	200	5.5	9.2		
Selenium	5.0	.38	.5	0.18	<5.0
Silicon	5.0	.38	.51		
Silver	3.0	.018	.051	-0.080	<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	0.49	<3.0

Associated samples MP7078: D32747-1

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

BLANK RESULTS SUMMARY  
Part 2 - Method Blanks

Login Number: D32747  
Account: XTOKRWR - XTO Energy  
Project: PCU T35X-2G

QC Batch ID: MP7078  
Matrix Type: SOLID

Methods: SW846 6010C  
Units: mg/kg

Prep Date:

Metal

(anr) Analyte not requested

## MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D32747  
 Account: XTOKRWR - XTO Energy  
 Project: PCU T35X-2G

QC Batch ID: MP7078  
 Matrix Type: SOLID

Methods: SW846 6010C  
 Units: mg/kg

Prep Date:

03/16/12

Metal	D32747-1 Original MS	Spikelot MPICPALL	% Rec	QC Limits
Aluminum				
Antimony				
Arsenic	anr			
Barium	1880	2000	207	58.0 (a) 75-125
Beryllium				
Boron				
Cadmium	0.18	44.5	51.8	85.6 75-125
Calcium				
Chromium	31.6	75.5	51.8	84.8 75-125
Cobalt				
Copper	15.5	63.7	51.8	93.1 75-125
Iron				
Lead	14.8	98.6	104	81.0 75-125
Lithium				
Magnesium				
Manganese				
Molybdenum				
Nickel	16.4	56.9	51.8	78.3 75-125
Phosphorus				
Potassium				
Selenium	0.96	86.5	104	82.6 75-125
Silicon				
Silver	0.26	19.4	20.7	92.5 75-125
Sodium				
Strontium				
Thallium				
Tin				
Titanium				
Uranium				
Vanadium				
Zinc	48.2	87.5	51.8	75.9 75-125

Associated samples MP7078: D32747-1

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

MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D32747  
Account: XTOKRWR - XTO Energy  
Project: PCU T35X-2G

QC Batch ID: MP7078  
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: D32747  
 Account: XTOKRWR - XTO Energy  
 Project: PCU T35X-2G

QC Batch ID: MP7078  
 Matrix Type: SOLID

Methods: SW846 6010C  
 Units: mg/kg

Prep Date:

03/16/12

Metal	D32747-1 Original	MSD	Spikelot MPICPALL	% Rec	MSD RPD	QC Limit
Aluminum						
Antimony						
Arsenic	anr					
Barium	1880	2000	213	56.3 (a)	0.0	20
Beryllium						
Boron						
Cadmium	0.18	45.3	53.3	84.6	1.8	20
Calcium						
Chromium	31.6	77.5	53.3	86.1	2.6	20
Cobalt						
Copper	15.5	65.1	53.3	93.0	2.2	20
Iron						
Lead	14.8	103	107	82.7	4.4	20
Lithium						
Magnesium						
Manganese						
Molybdenum						
Nickel	16.4	58.6	53.3	79.2	2.9	20
Phosphorus						
Potassium						
Selenium	0.96	87.6	107	81.3	1.3	20
Silicon						
Silver	0.26	19.8	21.3	91.6	2.0	20
Sodium						
Strontium						
Thallium						
Tin						
Titanium						
Uranium						
Vanadium						
Zinc	48.2	88.9	53.3	76.4	1.6	20

Associated samples MP7078: D32747-1

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

MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D32747  
Account: XTOKRWR - XTO Energy  
Project: PCU T35X-2G

QC Batch ID: MP7078  
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: D32747  
 Account: XTOKRWR - XTO Energy  
 Project: PCU T35X-2G

QC Batch ID: MP7078  
 Matrix Type: SOLID

Methods: SW846 6010C  
 Units: mg/kg

Prep Date:

03/16/12

Metal	BSP Result	Spikelot MPICPALL	% Rec	QC Limits
Aluminum				
Antimony				
Arsenic	anr			
Barium	193	200	96.5	80-120
Beryllium				
Boron				
Cadmium	47.5	50	95.0	80-120
Calcium				
Chromium	49.8	50	99.6	80-120
Cobalt				
Copper	49.5	50	99.0	80-120
Iron				
Lead	96.3	100	96.3	80-120
Lithium				
Magnesium				
Manganese				
Molybdenum				
Nickel	47.6	50	95.2	80-120
Phosphorus				
Potassium				
Selenium	94.2	100	94.2	80-120
Silicon				
Silver	21.0	20	105.0	80-120
Sodium				
Strontium				
Thallium				
Tin				
Titanium				
Uranium				
Vanadium				
Zinc	47.4	50	94.8	80-120

Associated samples MP7078: D32747-1

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

SPIKE BLANK AND LAB CONTROL SAMPLE SUMMARY

Login Number: D32747  
Account: XTOKWR - XTO Energy  
Project: PCU T35X-2G

QC Batch ID: MP7078  
Matrix Type: SOLID

Methods: SW846 6010C  
Units: mg/kg

Prep Date:

Metal

(anr) Analyte not requested

13.1.3  
**13**

## SERIAL DILUTION RESULTS SUMMARY

Login Number: D32747  
 Account: XTOKWR - XTO Energy  
 Project: PCU T35X-2G

QC Batch ID: MP7078  
 Matrix Type: SOLID

Methods: SW846 6010C  
 Units: ug/l

Prep Date: 03/16/12

Metal	D32747-1 Original	SDL 1:5	%DIF	QC Limits
Aluminum				
Antimony				
Arsenic	anr			
Barium	17500	19800	13.3*(a)	0-10
Beryllium				
Boron				
Cadmium	1.70	0.00	100.0(b)	0-10
Calcium				
Chromium	293	331	12.7*(a)	0-10
Cobalt				
Copper	144	141	2.6	0-10
Iron				
Lead	138	146	5.7	0-10
Lithium				
Magnesium				
Manganese				
Molybdenum				
Nickel	152	178	16.9*(a)	0-10
Phosphorus				
Potassium				
Selenium	8.90	20.5	130.3(b)	0-10
Silicon				
Silver	2.40	3.50	45.8 (b)	0-10
Sodium				
Strontium				
Thallium				
Tin				
Titanium				
Uranium				
Vanadium				
Zinc	448	541	20.7*(a)	0-10

Associated samples MP7078: D32747-1

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

SERIAL DILUTION RESULTS SUMMARY

Login Number: D32747  
Account: XTOKRWR - XTO Energy  
Project: PCU T35X-2G

QC Batch ID: MP7078  
Matrix Type: SOLID

Methods: SW846 6010C  
Units: ug/l

Prep Date:

Metal

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

BLANK RESULTS SUMMARY  
Part 2 - Method Blanks

Login Number: D32747  
Account: XTOKRWR - XTO Energy  
Project: PCU T35X-2G

QC Batch ID: MP7079  
Matrix Type: SOLID

Methods: SW846 6020A  
Units: mg/kg

Prep Date:

03/16/12

Metal	RL	IDL	MDL	MB raw	final
Aluminum	25	.14	1.3		
Antimony	0.20	.001	.012		
Arsenic	0.40	.049		0.056	<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 MP7079: D32747-1

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

13.2.1  
**13**

## MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D32747  
 Account: XTOKWR - XTO Energy  
 Project: PCU T35X-2G

QC Batch ID: MP7079  
 Matrix Type: SOLID

Methods: SW846 6020A  
 Units: mg/kg

Prep Date: 03/16/12

Metal	D32747-1 Original MS	Spikelot MPICPALL	% Rec	QC Limits
Aluminum				
Antimony				
Arsenic	3.6	111	104	103.8 75-125
Barium				
Beryllium				
Boron				
Cadmium				
Calcium				
Chromium				
Cobalt				
Copper				
Iron				
Lead				
Magnesium				
Manganese				
Molybdenum				
Nickel				
Phosphorus				
Potassium				
Selenium				
Silver				
Sodium				
Strontium				
Thallium				
Tin				
Titanium				
Uranium				
Vanadium				
Zinc				

Associated samples MP7079: D32747-1

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: D32747  
 Account: XTOKRWR - XTO Energy  
 Project: PCU T35X-2G

QC Batch ID: MP7079  
 Matrix Type: SOLID

Methods: SW846 6020A  
 Units: mg/kg

Prep Date:

03/16/12

Metal	D32747-1 Original	MSD	Spikelot MPICPALL	% Rec	MSD RPD	QC Limit
Aluminum						
Antimony						
Arsenic	3.6	116	107	105.4	4.4	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 MP7079: D32747-1

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: D32747  
 Account: XTOKRWR - XTO Energy  
 Project: PCU T35X-2G

QC Batch ID: MP7079  
 Matrix Type: SOLID

Methods: SW846 6020A  
 Units: mg/kg

Prep Date: 03/16/12

Metal	BSP Result	Spikelot MPICPALL	% Rec	QC Limits
Aluminum				
Antimony				
Arsenic	103	100	103.0	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 MP7079: D32747-1

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

## SERIAL DILUTION RESULTS SUMMARY

Login Number: D32747  
 Account: XTOKWR - XTO Energy  
 Project: PCU T35X-2G

QC Batch ID: MP7079  
 Matrix Type: SOLID

Methods: SW846 6020A  
 Units: ug/l

Prep Date: 03/16/12

Metal	D32747-1 Original	SDL 5:25	%DIF	QC Limits
Aluminum				
Antimony				
Arsenic	33.6	31.5	6.3	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 MP7079: D32747-1

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: D32747  
Account: XTOKRWR - XTO Energy  
Project: PCU T35X-2G

QC Batch ID: MP7098  
Matrix Type: SOLID

Methods: SW846 7471B  
Units: mg/kg

Prep Date: 03/20/12

Metal	RL	IDL	MDL	MB raw	final
Mercury	0.10	.0011	.013	0.0016	<0.10

Associated samples MP7098: D32747-1

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: D32747  
Account: XTOKRWR - XTO Energy  
Project: PCU T35X-2G

QC Batch ID: MP7098  
Matrix Type: SOLID

Methods: SW846 7471B  
Units: mg/kg

Prep Date: 03/20/12

Metal	D32726-1 Original MS	Spikelot HGWSR1	QC % Rec	QC Limits
Mercury	0.019	0.40	0.415	91.7 75-125

Associated samples MP7098: D32747-1

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: D32747  
Account: XTOKRWR - XTO Energy  
Project: PCU T35X-2G

QC Batch ID: MP7098  
Matrix Type: SOLID

Methods: SW846 7471B  
Units: mg/kg

Prep Date:

03/20/12

Metal	D32726-1 Original	MSD	Spikelot HGWSR1	MSD % Rec	RPD	QC Limit
Mercury	0.019	0.43	0.432	95.1	7.2	

Associated samples MP7098: D32747-1

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: D32747  
Account: XTOKRWR - XTO Energy  
Project: PCU T35X-2G

QC Batch ID: MP7098  
Matrix Type: SOLID

Methods: SW846 7471B  
Units: mg/kg

Prep Date: 03/20/12

Metal	BSP Result	Spikelot HGWSR1	QC % Rec	QC Limits
Mercury	0.39	0.4	97.5	80-120

Associated samples MP7098: D32747-1

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: D32747  
Account: XTOKRWR - XTO Energy  
Project: PCU T35X-2G

QC Batch ID: MP7102  
Matrix Type: AQUEOUS

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

Prep Date:

03/20/12

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	-10	<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	-9.0	<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	11.0	<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 MP7102: D32747-1A

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

BLANK RESULTS SUMMARY  
Part 2 - Method Blanks

Login Number: D32747  
Account: XTOKWR - XTO Energy  
Project: PCU T35X-2G

QC Batch ID: MP7102  
Matrix Type: AQUEOUS

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

Prep Date:

Metal

(anr) Analyte not requested

## MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D32747  
 Account: XTOKWR - XTO Energy  
 Project: PCU T35X-2G

QC Batch ID: MP7102  
 Matrix Type: AQUEOUS

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

Prep Date: 03/20/12

Metal	D32771-1A Original MS	Spikelot MPICPALL	% Rec	QC Limits
Aluminum				
Antimony				
Arsenic				
Barium				
Beryllium				
Boron				
Cadmium				
Calcium	152000	284000	125000	105.6    75-125
Chromium				
Cobalt				
Copper				
Iron				
Lead				
Lithium				
Magnesium	34700	163000	125000	102.6    75-125
Manganese				
Molybdenum				
Nickel				
Phosphorus				
Potassium				
Selenium				
Silicon				
Silver				
Sodium	128000	260000	125000	105.6    75-125
Strontium				
Thallium				
Tin				
Titanium				
Uranium				
Vanadium				
Zinc				

Associated samples MP7102: D32747-1A

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

MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D32747  
Account: XTOKWR - XTO Energy  
Project: PCU T35X-2G

QC Batch ID: MP7102  
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: D32747  
 Account: XTOKRWR - XTO Energy  
 Project: PCU T35X-2G

QC Batch ID: MP7102  
 Matrix Type: AQUEOUS

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

Prep Date: 03/20/12

Metal	D32771-1A Original MSD	Spikelot MPICPALL	MSD % Rec	MSD RPD	QC Limit
Aluminum					
Antimony					
Arsenic					
Barium					
Beryllium					
Boron					
Cadmium					
Calcium	152000	289000	125000	109.6	1.7
Chromium					
Cobalt					
Copper					
Iron					
Lead					
Lithium					
Magnesium	34700	164000	125000	103.4	0.6
Manganese					
Molybdenum					
Nickel					
Phosphorus					
Potassium					
Selenium					
Silicon					
Silver					
Sodium	128000	264000	125000	108.8	1.5
Strontium					
Thallium					
Tin					
Titanium					
Uranium					
Vanadium					
Zinc					

Associated samples MP7102: D32747-1A

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

MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D32747  
Account: XTOKWR - XTO Energy  
Project: PCU T35X-2G

QC Batch ID: MP7102  
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: D32747  
 Account: XTOKRWR - XTO Energy  
 Project: PCU T35X-2G

QC Batch ID: MP7102  
 Matrix Type: AQUEOUS

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

Prep Date: 03/20/12

Metal	BSP Result	Spikelot MPICPALL	% Rec	QC Limits
Aluminum				
Antimony				
Arsenic				
Barium				
Beryllium				
Boron				
Cadmium				
Calcium	128000	125000	102.4	80-120
Chromium				
Cobalt				
Copper				
Iron				
Lead				
Lithium				
Magnesium	126000	125000	100.8	80-120
Manganese				
Molybdenum				
Nickel				
Phosphorus				
Potassium				
Selenium				
Silicon				
Silver				
Sodium	129000	125000	103.2	80-120
Strontium				
Thallium				
Tin				
Titanium				
Uranium				
Vanadium				
Zinc				

Associated samples MP7102: D32747-1A

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

SPIKE BLANK AND LAB CONTROL SAMPLE SUMMARY

Login Number: D32747  
Account: XTOKWR - XTO Energy  
Project: PCU T35X-2G

QC Batch ID: MP7102  
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: D32747  
Account: XTOKWR - XTO Energy  
Project: PCU T35X-2G

Analyte	Batch ID	RL	MB Result	Units	Spike Amount	BSP Result	BSP %Recov	QC Limits
Chromium, Hexavalent	GP6737/GN14152	1.0	0.0	mg/kg	185	178	96.4	80-120%
Specific Conductivity	GP6755/GN14169	1.0	<1.0	umhos/cm	9967	9890	99.2	90-110%
pH	GN14113			su	8.00	7.96	99.5	99.3-100.7%

Associated Samples:

Batch GN14113: D32747-1

Batch GP6737: D32747-1

Batch GP6755: D32747-1

(\*) Outside of QC limits

DUPLICATE RESULTS SUMMARY  
GENERAL CHEMISTRY

Login Number: D32747  
Account: XTOKWR - XTO Energy  
Project: PCU T35X-2G

Analyte	Batch ID	QC Sample	Units	Original Result	DUP Result	RPD	QC Limits
Chromium, Hexavalent Redox Potential Vs H2	GP6737/GN14152 GN14122	D32651-1 D32747-1	mg/kg mv	0.0 377	0.0 361	0.0 4.3	0-20% 0-20%

Associated Samples:

Batch GN14122: D32747-1

Batch GP6737: D32747-1

(\*) Outside of QC limits

MATRIX SPIKE RESULTS SUMMARY  
GENERAL CHEMISTRY

Login Number: D32747  
Account: XTOKWR - XTO Energy  
Project: PCU T35X-2G

Analyte	Batch ID	QC Sample	Units	Original Result	Spike Amount	MS Result	%Rec	QC Limits
Chromium, Hexavalent	GP6737/GN14152	D32651-1	mg/kg	0.0	40	46.1	115.0	75-125%

Associated Samples:

Batch GP6737: D32747-1

(\*) Outside of QC limits

(N) Matrix Spike Rec. outside of QC limits

MATRIX SPIKE DUPLICATE RESULTS SUMMARY  
GENERAL CHEMISTRY

Login Number: D32747  
Account: XTOKWR - XTO Energy  
Project: PCU T35X-2G

Analyte	Batch ID	QC Sample	Units	Original Result	Spike Amount	MSD Result	RPD	QC Limit
Chromium, Hexavalent	GP6737/GN14152	D32651-1	mg/kg	0.0	40	44.3	4.1	

Associated Samples:

Batch GP6737: D32747-1

(\*) Outside of QC limits

(N) Matrix Spike Rec. outside of QC limits