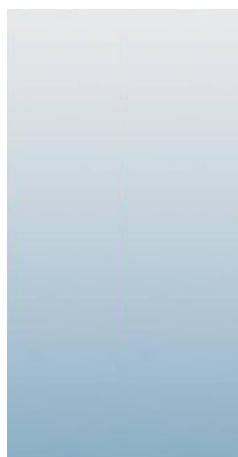




01/11/13



## Technical Report for

**XTO Energy**

**XTO Love Ranch 8**

**1108-07A**

**Accutest Job Number: D42316**

**Sampling Date: 01/03/13**

### Report to:

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

**Total number of pages in report: 169**



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.

# Table of Contents

-1-

<b>Section 1: Sample Summary .....</b>	<b>4</b>	
<b>Section 2: Case Narrative/Conformance Summary .....</b>	<b>5</b>	
<b>Section 3: Summary of Hits .....</b>	<b>8</b>	
<b>Section 4: Sample Results .....</b>	<b>9</b>	
<b>4.1: D42316-1: RP POST SOLIDIFICATION .....</b>	<b>10</b>	
<b>4.2: D42316-1A: RP POST SOLIDIFICATION .....</b>	<b>16</b>	
<b>Section 5: Misc. Forms .....</b>	<b>18</b>	
<b>5.1: Chain of Custody .....</b>	<b>19</b>	
<b>Section 6: GC/MS Volatiles - QC Data Summaries .....</b>	<b>21</b>	
<b>6.1: Method Blank Summary .....</b>	<b>22</b>	
<b>6.2: Blank Spike Summary .....</b>	<b>23</b>	
<b>6.3: Matrix Spike/Matrix Spike Duplicate Summary .....</b>	<b>24</b>	
<b>Section 7: GC/MS Volatiles - Raw Data .....</b>	<b>25</b>	
<b>7.1: Samples .....</b>	<b>26</b>	
<b>7.2: Method Blanks .....</b>	<b>41</b>	
<b>Section 8: GC/MS Semi-volatiles - QC Data Summaries .....</b>	<b>53</b>	
<b>8.1: Method Blank Summary .....</b>	<b>54</b>	
<b>8.2: Blank Spike Summary .....</b>	<b>55</b>	
<b>8.3: Matrix Spike/Matrix Spike Duplicate Summary .....</b>	<b>56</b>	
<b>Section 9: GC/MS Semi-volatiles - Raw Data .....</b>	<b>57</b>	
<b>9.1: Samples .....</b>	<b>58</b>	
<b>9.2: Method Blanks .....</b>	<b>92</b>	
<b>Section 10: GC Volatiles - QC Data Summaries .....</b>	<b>109</b>	
<b>10.1: Method Blank Summary .....</b>	<b>110</b>	
<b>10.2: Blank Spike Summary .....</b>	<b>111</b>	
<b>10.3: Matrix Spike/Matrix Spike Duplicate Summary .....</b>	<b>112</b>	
<b>Section 11: GC Volatiles - Raw Data .....</b>	<b>113</b>	
<b>11.1: Samples .....</b>	<b>114</b>	
<b>11.2: Method Blanks .....</b>	<b>119</b>	
<b>Section 12: GC Semi-volatiles - QC Data Summaries .....</b>	<b>124</b>	
<b>12.1: Method Blank Summary .....</b>	<b>125</b>	
<b>12.2: Blank Spike Summary .....</b>	<b>126</b>	
<b>12.3: Matrix Spike/Matrix Spike Duplicate Summary .....</b>	<b>127</b>	
<b>Section 13: GC Semi-volatiles - Raw Data .....</b>	<b>128</b>	
<b>13.1: Samples .....</b>	<b>129</b>	
<b>13.2: Method Blanks .....</b>	<b>132</b>	
<b>Section 14: Metals Analysis - QC Data Summaries .....</b>	<b>135</b>	
<b>14.1: Prep QC MP9199: Ba,Cd,Cr,Cu,Pb,Ni,Se,Ag,Zn .....</b>	<b>136</b>	
<b>14.2: Prep QC MP9200: As .....</b>	<b>146</b>	
<b>14.3: Prep QC MP9202: Hg .....</b>	<b>151</b>	
<b>14.4: Prep QC MP9206: Ca,Mg,Na,Sodium Adsorption Ratio .....</b>	<b>155</b>	
<b>Section 15: General Chemistry - QC Data Summaries .....</b>	<b>165</b>	

# Table of Contents

-2-

<b>15.1:</b> Method Blank and Spike Results Summary .....	166
<b>15.2:</b> Duplicate Results Summary .....	167
<b>15.3:</b> Matrix Spike Results Summary .....	168
<b>15.4:</b> Matrix Spike Duplicate Results Summary .....	169

1  
2  
3  
4  
5  
6  
7  
8  
9  
10  
11  
12  
13  
14  
15



## Sample Summary

XTO Energy

**Job No:** D42316

XTO Love Ranch 8

Project No: 1108-07A

Sample Number	Collected Date	Time By	Matrix Received	Code Type	Client Sample ID
D42316-1	01/03/13	10:00 DS	01/05/13	SO Soil	RP POST SOLIDIFICATION
D42316-1A	01/03/13	10:00 DS	01/05/13	SO Soil	RP POST SOLIDIFICATION

---

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



## CASE NARRATIVE / CONFORMANCE SUMMARY

**Client:** XTO Energy

**Job No** D42316

**Site:** XTO Love Ranch 8

**Report Dat** 1/11/2013 1:28:14 PM

On 01/05/2013, 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 D42316 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> V3V1321
------------------	--------------------------

- All samples were analyzed within the recommended method holding time.
- All method blanks for this batch meet method specific criteria.
- Sample(s) D42318-1MS, D42318-1MSD were used as the QC samples indicated.

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

<b>Matrix</b> SO	<b>Batch ID:</b> OP7200
------------------	-------------------------

- All samples were extracted and analyzed within the recommended method holding time.
- Sample(s) D42316-1MS, D42316-1MSD were used as the QC samples indicated.
- All method blanks for this batch meet method specific criteria.
- The matrix spike (MS) recovery(s) of Fluorene are outside control limits. Outside control limits due to high level in sample relative to spike amount.
- The matrix spike duplicate (MSD) recovery(s) of Acenaphthene are outside control limits. Variability of recovery may be due to sample matrix/homogeneity.
- The matrix spike (MS) recovery(s) of Naphthalene are outside control limits. Outside control limits due to high level in sample relative to spike amount.
- The RPD(s) for the MS and MSD recoveries of Acenaphthene are outside control limits for sample OP7200-MSD. Variability of recovery may be due to sample matrix/homogeneity.
- Sample(s) OP7200-MB have surrogates outside control limits. Probable cause due to matrix interference.
- OP7200-MB for Terphenyl-d14: Outside of control limits. Since the bias is high and the method blank is ND for target analytes, no further action is required.

### Volatiles by GC By Method SW846 8015B

<b>Matrix</b> SO	<b>Batch ID:</b> GGB1042
------------------	--------------------------

- All samples were analyzed within the recommended method holding time.
- Sample(s) D42317-1MS, D42317-1MSD were used as the QC samples indicated.
- All method blanks for this batch meet method specific criteria.

## Extractables by GC By Method SW846-8015B

**Matrix** SO

**Batch ID:** OP7201

- All samples were extracted and analyzed within the recommended method holding time.
- All method blanks for this batch meet method specific criteria.
- Sample(s) D42316-1MS, D42316-1MSD were used as the QC samples indicated.
- The matrix spike (MS) recovery(s) of TPH-DRO (C10-C28) are outside control limits. Outside control limits due to high level in sample relative to spike amount.

## Metals By Method SW846 6010C

**Matrix** AQ

**Batch ID:** MP9206

- All samples were digested and analyzed within the recommended method holding time.
- All method blanks for this batch meet method specific criteria.
- Sample(s) D42292-1AMS, D42292-1AMSD, D42292-1ASDL were used as the QC samples for the metals analysis.

**Matrix** SO

**Batch ID:** MP9199

- All samples were digested and analyzed within the recommended method holding time.
- All method blanks for this batch meet method specific criteria.
- Sample(s) D42292-1MS, D42292-1MSD, D42292-1SDL were used as the QC samples for the metals analysis.
- The serial dilution RPD(s) for Cadmium, Copper, Lead, Barium, Zinc are outside control limits for sample MP9199-SD1. Percent difference acceptable due to low initial sample concentration (< 50 times IDL).
- MP9199-SD1 for Zinc: Serial dilution indicates possible matrix interference.
- MP9199-SD1 for Barium: Serial dilution indicates possible matrix interference.

## Metals By Method SW846 6020A

**Matrix** SO

**Batch ID:** MP9200

- All samples were digested and analyzed within the recommended method holding time.
- All method blanks for this batch meet method specific criteria.
- Sample(s) D42292-1MS, D42292-1MSD, D42292-1SDL were used as the QC samples for the metals analysis.

## Metals By Method SW846 7471B

**Matrix** SO

**Batch ID:** MP9202

- All samples were digested and analyzed within the recommended method holding time.
- All method blanks for this batch meet method specific criteria.
- Sample(s) D42292-1MS, D42292-1MSD were used as the QC samples for the metals analysis.

## Wet Chemistry By Method ASTM D1498-76M

**Matrix** SO

**Batch ID:** GN18332

- Sample(s) D42316-1DUP were used as the QC samples for the Redox Potential Vs H<sub>2</sub> analysis.

## Wet Chemistry By Method SM19 2540B M

**Matrix** SO

**Batch ID:** GN18327

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

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

**Matrix** SO

**Batch ID:** GP9044

- All samples were prepared and analyzed within the recommended method holding time.
- All method blanks for this batch meet method specific criteria.
- Sample(s) D42316-1DUP, D42316-1MS, D42316-1MSD were used as the QC samples for the Chromium, Hexavalent analysis.
- D42316-1 for Chromium, Hexavalent: Dilution required due to matrix interference.

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

**Matrix** SO

**Batch ID:** R15651

- The data for SW846 3060A/7196A M meets quality control requirements.
- D42316-1 for Chromium, Trivalent: Calculated as: (Chromium) - (Chromium, Hexavalent)

## Wet Chemistry By Method SW846 9045D

**Matrix** SO

**Batch ID:** GN18331

- The following samples were run outside of holding time for method SW846 9045D: D42316-1

## Wet Chemistry By Method USDA HANDBOOK 60

**Matrix** SO

**Batch ID:** MP9206

- D42316-1A for Sodium Adsorption Ratio: Calculated as:  $(\text{Na meq/L}) / \sqrt{[(\text{Ca meq/L}) + (\text{Mg meq/L})] / 2}$

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

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

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

## Summary of Hits

Page 1 of 1

Job Number: D42316  
Account: XTO Energy  
Project: XTO Love Ranch 8  
Collected: 01/03/13

3

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

### D42316-1 RP POST SOLIDIFICATION

Toluene	0.378	0.22	0.11	mg/kg	SW846 8260B
Ethylbenzene	0.336	0.22	0.041	mg/kg	SW846 8260B
Xylene (total)	9.32	0.43	0.22	mg/kg	SW846 8260B
Fluorene	0.437	0.013	0.0068	mg/kg	SW846 8270C BY SIM
Naphthalene	1.19	0.074	0.065	mg/kg	SW846 8270C BY SIM
Pyrene	0.0800	0.013	0.0068	mg/kg	SW846 8270C BY SIM
TPH-GRO (C6-C10)	225	22	11	mg/kg	SW846 8015B
TPH-DRO (C10-C28)	5850	110	63	mg/kg	SW846-8015B
Arsenic	8.8	0.15		mg/kg	SW846 6020A
Barium	8260	7.7		mg/kg	SW846 6010C
Chromium	20.0	1.5		mg/kg	SW846 6010C
Copper	25.2	1.5		mg/kg	SW846 6010C
Lead	8.9	7.7		mg/kg	SW846 6010C
Nickel	86.2	4.6		mg/kg	SW846 6010C
Zinc	38.5	4.6		mg/kg	SW846 6010C
Specific Conductivity	12400	1.0		umhos/cm	SM 2510B-2011 MOD
Chromium, Trivalent <sup>a</sup>	20.0	6.5		mg/kg	SW846 3060A/7196A M
Redox Potential Vs H2	36.3			mv	ASTM D1498-76M
pH	11.33			su	SW846 9045D

### D42316-1A RP POST SOLIDIFICATION

Calcium	623	2.0	mg/l	SW846 6010C
Sodium	1470	2.0	mg/l	SW846 6010C
Sodium Adsorption Ratio <sup>b</sup>	16.2		ratio	USDA HANDBOOK 60

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

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



4

## Sample Results

---

### Report of Analysis

---

Accutest Laboratories

**Report of Analysis**

Page 1 of 1

**Client Sample ID:** RP POST SOLIDIFICATION**Lab Sample ID:** D42316-1**Date Sampled:** 01/03/13**Matrix:** SO - Soil**Date Received:** 01/05/13**Method:** SW846 8260B**Percent Solids:** 63.3**Project:** XTO Love Ranch 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	3V22448.D	1	01/08/13	BD	n/a	n/a	V3V1321
Run #2							

	<b>Initial Weight</b>	<b>Final Volume</b>	<b>Methanol Aliquot</b>
Run #1	5.02 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	ND	0.11	0.054	mg/kg	
108-88-3	Toluene	0.378	0.22	0.11	mg/kg	
100-41-4	Ethylbenzene	0.336	0.22	0.041	mg/kg	
1330-20-7	Xylene (total)	9.32	0.43	0.22	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	90%		64-130%
460-00-4	4-Bromofluorobenzene	120%		62-131%
17060-07-0	1,2-Dichloroethane-D4	101%		70-130%

ND = Not detected MDL - Method Detection Limit

J = Indicates an estimated value

RL = Reporting Limit

B = Indicates analyte found in associated method blank

E = Indicates value exceeds calibration range

N = Indicates presumptive evidence of a compound

Accutest Laboratories

**Report of Analysis**

Page 1 of 1

<b>Client Sample ID:</b>	RP POST SOLIDIFICATION	<b>Date Sampled:</b>	01/03/13
<b>Lab Sample ID:</b>	D42316-1	<b>Date Received:</b>	01/05/13
<b>Matrix:</b>	SO - Soil	<b>Percent Solids:</b>	63.3
<b>Method:</b>	SW846 8270C BY SIM SW846 3546		
<b>Project:</b>	XTO Love Ranch 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	3G12922.D	1	01/10/13	DC	01/09/13	OP7200	E3G618
Run #2	3G12912.D	4	01/10/13	DC	01/09/13	OP7200	E3G618

	<b>Initial Weight</b>	<b>Final Volume</b>
Run #1	30.0 g	1.0 ml
Run #2	30.0 g	1.0 ml

**COGCC Table 910-1 PAH List**

<b>CAS No.</b>	<b>Compound</b>	<b>Result</b>	<b>RL</b>	<b>MDL</b>	<b>Units</b>	<b>Q</b>
83-32-9	Acenaphthene	ND	0.013	0.0068	mg/kg	
120-12-7	Anthracene	ND	0.013	0.0068	mg/kg	
56-55-3	Benzo(a)anthracene	ND	0.013	0.0068	mg/kg	
205-99-2	Benzo(b)fluoranthene	ND	0.013	0.0068	mg/kg	
207-08-9	Benzo(k)fluoranthene	ND	0.013	0.0068	mg/kg	
50-32-8	Benzo(a)pyrene	ND	0.013	0.0068	mg/kg	
218-01-9	Chrysene	ND	0.013	0.0068	mg/kg	
53-70-3	Dibenz(a,h)anthracene	ND	0.013	0.0068	mg/kg	
206-44-0	Fluoranthene	ND	0.013	0.0068	mg/kg	
86-73-7	Fluorene	0.437	0.013	0.0068	mg/kg	
193-39-5	Indeno(1,2,3-cd)pyrene	ND	0.013	0.0068	mg/kg	
91-20-3	Naphthalene	1.19 <sup>a</sup>	0.074	0.065	mg/kg	
129-00-0	Pyrene	0.0800	0.013	0.0068	mg/kg	

<b>CAS No.</b>	<b>Surrogate Recoveries</b>	<b>Run# 1</b>	<b>Run# 2</b>	<b>Limits</b>
4165-60-0	Nitrobenzene-d5	84%	42%	10-159%
321-60-8	2-Fluorobiphenyl	50%	48%	19-131%
1718-51-0	Terphenyl-d14	73%	107%	18-150%

(a) Result is from Run# 2

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**

Page 1 of 1

**Client Sample ID:** RP POST SOLIDIFICATION**Lab Sample ID:** D42316-1**Date Sampled:** 01/03/13**Matrix:** SO - Soil**Date Received:** 01/05/13**Method:** SW846 8015B**Percent Solids:** 63.3**Project:** XTO Love Ranch 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	GB19081.D	1	01/07/13	SK	n/a	n/a	GGB1042
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)	225	22	11	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	123%			60-140%	

ND = Not detected MDL - Method Detection Limit

J = Indicates an estimated value

RL = Reporting Limit

B = Indicates analyte found in associated method blank

E = Indicates value exceeds calibration range

N = Indicates presumptive evidence of a compound

4.1

4

Accutest Laboratories

**Report of Analysis**

Page 1 of 1

**Client Sample ID:** RP POST SOLIDIFICATION**Lab Sample ID:** D42316-1**Date Sampled:** 01/03/13**Matrix:** SO - Soil**Date Received:** 01/05/13**Method:** SW846-8015B SW846 3546**Percent Solids:** 63.3**Project:** XTO Love Ranch 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	FD21072.D	10	01/10/13	AV	01/09/13	OP7201	GFD1060
Run #2							

	<b>Initial Weight</b>	<b>Final Volume</b>
Run #1	30.1 g	1.0 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)	5850	110	63	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	69%		35-130%		

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

4.1

4

**Report of Analysis**

Page 1 of 1

<b>Client Sample ID:</b>	RP POST SOLIDIFICATION	<b>Date Sampled:</b>	01/03/13
<b>Lab Sample ID:</b>	D42316-1	<b>Date Received:</b>	01/05/13
<b>Matrix:</b>	SO - Soil	<b>Percent Solids:</b>	63.3
<b>Project:</b>	XTO Love Ranch 8		

**Metals Analysis**

Analyte	Result	RL	Units	DF	Prep	Analyzed By	Method	Prep Method
Arsenic	8.8	0.15	mg/kg	5	01/07/13	01/08/13	JB	SW846 6020A <sup>1</sup>
Barium	8260	7.7	mg/kg	5	01/07/13	01/09/13	JB	SW846 6010C <sup>3</sup>
Cadmium	< 1.5	1.5	mg/kg	1	01/07/13	01/08/13	JB	SW846 6010C <sup>3</sup>
Chromium	20.0	1.5	mg/kg	1	01/07/13	01/08/13	JB	SW846 6010C <sup>3</sup>
Copper	25.2	1.5	mg/kg	1	01/07/13	01/08/13	JB	SW846 6010C <sup>3</sup>
Lead	8.9	7.7	mg/kg	1	01/07/13	01/08/13	JB	SW846 6010C <sup>3</sup>
Mercury	< 0.12	0.12	mg/kg	1	01/08/13	01/08/13	JM	SW846 7471B <sup>2</sup>
Nickel	86.2	4.6	mg/kg	1	01/07/13	01/08/13	JB	SW846 6010C <sup>3</sup>
Selenium	< 7.7	7.7	mg/kg	1	01/07/13	01/08/13	JB	SW846 6010C <sup>3</sup>
Silver	< 4.6	4.6	mg/kg	1	01/07/13	01/08/13	JB	SW846 6010C <sup>3</sup>
Zinc	38.5	4.6	mg/kg	1	01/07/13	01/08/13	JB	SW846 6010C <sup>3</sup>

- (1) Instrument QC Batch: MA3150
- (2) Instrument QC Batch: MA3152
- (3) Instrument QC Batch: MA3153
- (4) Prep QC Batch: MP9199
- (5) Prep QC Batch: MP9200
- (6) Prep QC Batch: MP9202

RL = Reporting Limit

**Report of Analysis**

Page 1 of 1

**Client Sample ID:** RP POST SOLIDIFICATION**Lab Sample ID:** D42316-1**Matrix:** SO - Soil**Date Sampled:** 01/03/13**Date Received:** 01/05/13**Percent Solids:** 63.3**Project:** XTO Love Ranch 8**General Chemistry**

Analyte	Result	RL	Units	DF	Analyzed	By	Method
<b>prep: DEPT.OF AG, BOOK N9</b>							
Specific Conductivity	12400	1.0	umhos/cm	1	01/09/13	JD	SM 2510B-2011 MOD
Chromium, Hexavalent <sup>a</sup>	< 5.0	5.0	mg/kg	5	01/07/13	KB	SW846 3060A/7196A
Chromium, Trivalent <sup>b</sup>	20.0	6.5	mg/kg	1	01/08/13 21:55	JB	SW846 3060A/7196A M
Redox Potential Vs H2	36.3		mv	1	01/07/13	CT	ASTM D1498-76M
Solids, Percent	63.3		%	1	01/07/13	SWT	SM19 2540B M
pH	11.33		su	1	01/07/13 13:00	CT	SW846 9045D

(a) Dilution required due to matrix interference.

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

RL = Reporting Limit

**Report of Analysis**

Page 1 of 1

<b>Client Sample ID:</b>	RP POST SOLIDIFICATION	<b>Date Sampled:</b>	01/03/13
<b>Lab Sample ID:</b>	D42316-1A	<b>Date Received:</b>	01/05/13
<b>Matrix:</b>	SO - Soil	<b>Percent Solids:</b>	63.3
<b>Project:</b>	XTO Love Ranch 8		

**SAR Metals Analysis**

Analyte	Result	RL	Units	DF	Prep	Analyzed By	Method	Prep Method
Calcium	623	2.0	mg/l	1	01/08/13	01/08/13 JB	SW846 6010C <sup>1</sup>	SW846 3010A/M <sup>2</sup>
Magnesium	< 1.0	1.0	mg/l	1	01/08/13	01/08/13 JB	SW846 6010C <sup>1</sup>	SW846 3010A/M <sup>2</sup>
Sodium	1470	2.0	mg/l	1	01/08/13	01/08/13 JB	SW846 6010C <sup>1</sup>	SW846 3010A/M <sup>2</sup>

(1) Instrument QC Batch: MA3153

(2) Prep QC Batch: MP9206

RL = Reporting Limit

**Report of Analysis**

Page 1 of 1

**Client Sample ID:** RP POST SOLIDIFICATION**Lab Sample ID:** D42316-1A**Matrix:** SO - Soil**Date Sampled:** 01/03/13**Date Received:** 01/05/13**Percent Solids:** 63.3**Project:** XTO Love Ranch 8**General Chemistry**

Analyte	Result	RL	Units	DF	Analyzed	By	Method
Sodium Adsorption Ratio <sup>a</sup>	16.2		ratio	1	01/08/13 18:13	JB	USDA HANDBOOK 60

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

RL = Reporting Limit

4.2

4



## Misc. Forms

---

5

### Custody Documents and Other Forms

---

Includes the following where applicable:

- Chain of Custody



## CHAIN OF CUSTODY

PAGE 1 OF 1

4036 Youngfield Street, Wheat Ridge, CO 80033  
TEL: 303-425-6021 FAX: 303-425-5854  
www.accutest.com

FED-EX Tracking #	Bottle Order Control #
Accutest Quote #	Accutest Job # D42316

Client / Reporting Information		Project Information		Requested Analysis ( see TEST CODE sheet)												Matrix Codes			
Company Name <b>KRW Consulting</b>	Project Name: <b>XTO LOVE RANCH 8</b>	Street Address <b>8000 West 14th Street; Suite 200</b>	City <b>Lakewood, CO 80214</b>	Billing Information ( If different from Report to )												DW - Drinking Water GW - Ground Water WW - Water SW - Surface Water SO - Soil SL - Sludge SE - Sediment OI - Oil LIQ - Other Liquid AIR - Air SOL - Other Solid WP - Wipe FB - Field Blank EB - Equipment Blank RB - Rinse Blank TB - Trip Blank			
Project Contact <b>Dwayne Knudson</b>	Project # <b>1108-07A</b>	Client Purchase Order # <b>970-488-1098</b>	City <b>Rifle, CO 81650</b>	Company Name <b>XTO Energy</b>															
Phone # <b>970-488-1098</b>	Project Manager <b>DAVID SANDERS</b>	Attention: <b>Joe Hess</b>	City <b>Rifle, CO 81650</b>	Street Address <b>21469 CR 5</b>															
Sampler(s) Name(s) <b>DAVID SANDERS</b>	Project Manager <b>Joe Hess</b>	Attention: <b>Jessica Dooling</b>	City <b>Rifle, CO 81650</b>	Number of preserved Bottles															
Accutest Sample #	Field ID / Point of Collection <b>RP Post Solidification</b>	Collection MEOH/DI Vial # <b>1-3-12</b>	Date <b>10:00</b>	Time <b>DS</b>	Sampled by <b>SO 5</b>	Matrix <b>HCl</b>	# of bottles <b>1</b>	HCl <input type="checkbox"/>	NaOH <input type="checkbox"/>	HNO3 <input type="checkbox"/>	H2SO4 <input type="checkbox"/>	None <input type="checkbox"/>	DI Water <input type="checkbox"/>	MECH <input type="checkbox"/>	ENCORE <input type="checkbox"/>	Blanks <input type="checkbox"/>	T-910	X	OJ
Turnaround Time ( Business days )																Data Deliverable Information	Comments / Special Instructions		
<input type="checkbox"/> Std. 10 Business Days <input checked="" type="checkbox"/> Std. 5 Business Days (By contract only) <input type="checkbox"/> 3 Day Emergency <input type="checkbox"/> 2 Day Emergency <input type="checkbox"/> 1 Day Emergency <input type="checkbox"/> Emergency & Rush T/A data available VIA Lablink																<input type="checkbox"/> Commercial "A" (Level 1) <input type="checkbox"/> Commercial "B" (Level 2) <input type="checkbox"/> COMMNB <input type="checkbox"/> COMMNB+ <input type="checkbox"/>	<input type="checkbox"/> State Forms Required <input type="checkbox"/> Send Forms to State <input type="checkbox"/> Report by Fax <input checked="" type="checkbox"/> Report by PDF ONLY <input type="checkbox"/> EDD Format	Please email to: <b>KRW Piceance Team</b>	
<b>Sample Custody must be documented below each time samples change possession, including courier delivery.</b>																			
Relinquished by Sampler: <b>1 Lori Albrecht</b>	Date Time: <b>1/14/13 12:00</b>	Received By: <b>KR Service Center</b>	Relinquished By: <b>2</b>	Date Time:	Received By:														
Relinquished by Sampler: <b>3</b>	Date Time:	Received By: <b>3</b>	Relinquished By: <b>4</b>	Date Time: <b>1/15/13 11:05</b>	Received By:	<b>4</b>													
Relinquished by: <b>5</b>	Date Time:	Received By: <b>5</b>	Custody Seal # <b>EX</b>	<input type="checkbox"/> Handled <input type="checkbox"/> Not Handled	Preservative chars applicable	<b>2</b>	On Ice <b>40°C</b>												

5.1

**D42316: Chain of Custody**  
**Page 1 of 2**



## Accutest Laboratories Sample Receipt Summary

Accutest Job Number: D42316

Client: KRW

Immediate Client Services Action Required: No

Date / Time Received: 1/5/2013 11:05:00 AM

No. Coolers:

1

Client Service Action Required at Login: No

Project: XTO LOVE RANCH 8

Airbill #'s: FX

**Cooler Security**Y or NY or N

1. Custody Seals Present:        3. COC Present:    
 2. Custody Seals Intact:        4. Smpl Dates/Time OK

**Cooler Temperature**Y or N

1. Temp criteria achieved:    
 2. Cooler temp verification: Infared gun  
 3. Cooler media: Ice (bag)

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

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

**Sample Integrity - Documentation**Y or N

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

**Sample Integrity - Condition**Y or N

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

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

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

Comments

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

5.1

5

**D42316: Chain of Custody****Page 2 of 2**



## GC/MS Volatiles

---

### QC Data Summaries

---

Includes the following where applicable:

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



**Method Blank Summary**

**Job Number:** D42316  
**Account:** XTOKWR XTO Energy  
**Project:** XTO Love Ranch 8

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
V3V1321-MB	3V22443.D	1	01/08/13	BD	n/a	n/a	V3V1321

**The QC reported here applies to the following samples:****Method:** SW846 8260B

D42316-1

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

CAS No.	Surrogate Recoveries	Limits
2037-26-5	Toluene-D8	85% 64-130%
460-00-4	4-Bromofluorobenzene	95% 62-131%
17060-07-0	1,2-Dichloroethane-D4	109% 70-130%

**Blank Spike Summary**

**Job Number:** D42316  
**Account:** XTOKWR XTO Energy  
**Project:** XTO Love Ranch 8

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
V3V1321-BS	3V22444.D	1	01/08/13	BD	n/a	n/a	V3V1321

**The QC reported here applies to the following samples:****Method:** SW846 8260B

D42316-1

CAS No.	Compound	Spike ug/kg	BSP ug/kg	BSP %	Limits
71-43-2	Benzene	50	52.4	105	70-130
100-41-4	Ethylbenzene	50	50.4	101	70-130
108-88-3	Toluene	50	49.0	98	70-130
1330-20-7	Xylene (total)	150	155	103	70-130

CAS No.	Surrogate Recoveries	BSP	Limits
2037-26-5	Toluene-D8	94%	64-130%
460-00-4	4-Bromofluorobenzene	104%	62-131%
17060-07-0	1,2-Dichloroethane-D4	102%	70-130%

\* = Outside of Control Limits.

# Matrix Spike/Matrix Spike Duplicate Summary

Page 1 of 1

Job Number: D42316

Account: XTOKWR XTO Energy

Project: XTO Love Ranch 8

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
D42318-1MS	3V22446.D	1	01/08/13	BD	n/a	n/a	V3V1321
D42318-1MSD	3V22447.D	1	01/08/13	BD	n/a	n/a	V3V1321
D42318-1	3V22445.D	1	01/08/13	BD	n/a	n/a	V3V1321

The QC reported here applies to the following samples:

Method: SW846 8260B

D42316-1

CAS No.	Compound	D42318-1		Spike	MS	MS	MSD	MSD	Limits	
		ug/kg	Q	ug/kg	ug/kg	%	ug/kg	%	RPD	Rec/RPD
71-43-2	Benzene	195		3310	3600	103	3540	101	2	64-139/30
100-41-4	Ethylbenzene	65.4	J	3310	3370	100	3310	98	2	68-136/30
108-88-3	Toluene	385		3310	3330	89	3330	89	0	60-130/30
1330-20-7	Xylene (total)	343		9920	10500	102	10300	100	2	58-142/30

CAS No.	Surrogate Recoveries	MS	MSD	D42318-1	Limits
2037-26-5	Toluene-D8	87%	88%	86%	64-130%
460-00-4	4-Bromofluorobenzene	108%	107%	99%	62-131%
17060-07-0	1,2-Dichloroethane-D4	98%	97%	106%	70-130%

\* = Outside of Control Limits.



GC/MS Volatiles

---

Raw Data

---

7

## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\V3010813.S\  
 Data File : 3V22448.D  
 Acq On : 8 Jan 2013 1:25 pm  
 Operator : BRETD  
 Sample : D42316-1  
 Misc : MS5203,V3V1321,5.022,,100,5,1  
 ALS Vial : 8 Sample Multiplier: 1

Quant Time: Jan 09 09:15:40 2013  
 Quant Method : C:\msdchem\1\METHODS\V3AP1299TVH1299SOIL.M  
 Quant Title : 8260  
 QLast Update : Thu Jan 03 11:40:16 2013  
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
2) Pentafluorobenzene	11.860	168	304850	50.00	ug/l	0.00
35) 1,4-Difluorobenzene	12.656	114	490467	50.00	ug/l	0.00
53) Chlorobenzene-d5	15.295	117	611188	50.00	ug/l	0.00
74) 1,4-Dichlorobenzene-d4	17.284	152	392865	50.00	ug/l	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
33) 1,2-Dichloroethane-d4	12.251	102	34391	50.41	ug/l	0.00
Spiked Amount 50.000	Range 70 - 130		Recovery	=	100.82%	
61) Toluene-d8	14.051	98	661180	44.93	ug/l	0.00
Spiked Amount 50.000	Range 70 - 130		Recovery	=	89.86%	
69) 4-Bromofluorobenzene	16.245	95	379777	59.91	ug/l	0.00
Spiked Amount 50.000	Range 70 - 130		Recovery	=	119.82%	

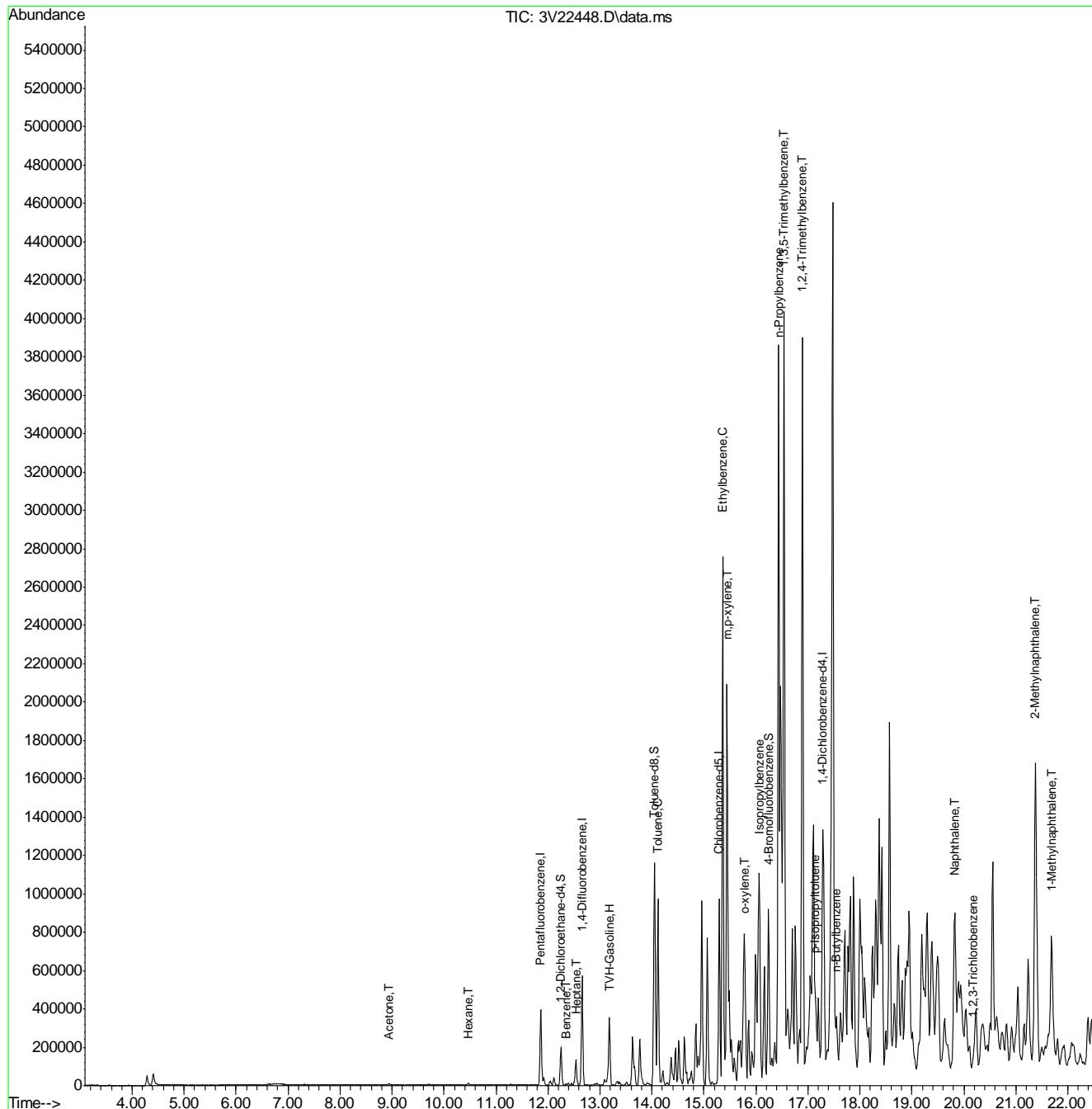
Target Compounds	R.T.	TIC	Qvalue
1) TVH-Gasoline	13.200	60115550m	2048.51 ug/l
15) Acetone	8.931	43	13321 6.95 ug/l 95
41) Hexane	10.465	57	4893 0.85 ug/l 100
43) Heptane	12.540	43	55500 8.43 ug/l 94
50) Benzene	12.348	78	6269 0.46 ug/l 100
62) Toluene	14.112	92	46727 3.51 ug/l 96
66) Ethylbenzene	15.366	91	69483 3.12 ug/l 98
68) Isopropylbenzene	16.078	105	65517 2.95 ug/l 97
72) m,p-xylene	15.446	106	683541 72.86 ug/l 98
73) o-xylene	15.796	106	124940 13.71 ug/l 98
77) n-Propylbenzene	16.425	91	251640 9.31 ug/l 95
80) 1,3,5-Trimethylbenzene	16.540	105	2232151 110.37 ug/l 93
82) 1,2,4-Trimethylbenzene	16.896	105	2462767 119.39 ug/l 93
86) p-Isopropyltoluene	17.153	119	166175 7.21 ug/l 99
88) n-Butylbenzene	17.541	91	103887 5.56 ug/l # 79
91) Naphthalene	19.841	128	621381 30.16 ug/l 100
93) 1,2,3-Trichlorobenzene	20.162	180	12733 1.55 ug/l 96
94) 2-Methylnaphthalene	21.377	142	1128845 123.72 ug/l 96
95) 1-Methylnaphthalene	21.685	142	403057 46.58 ug/l 95

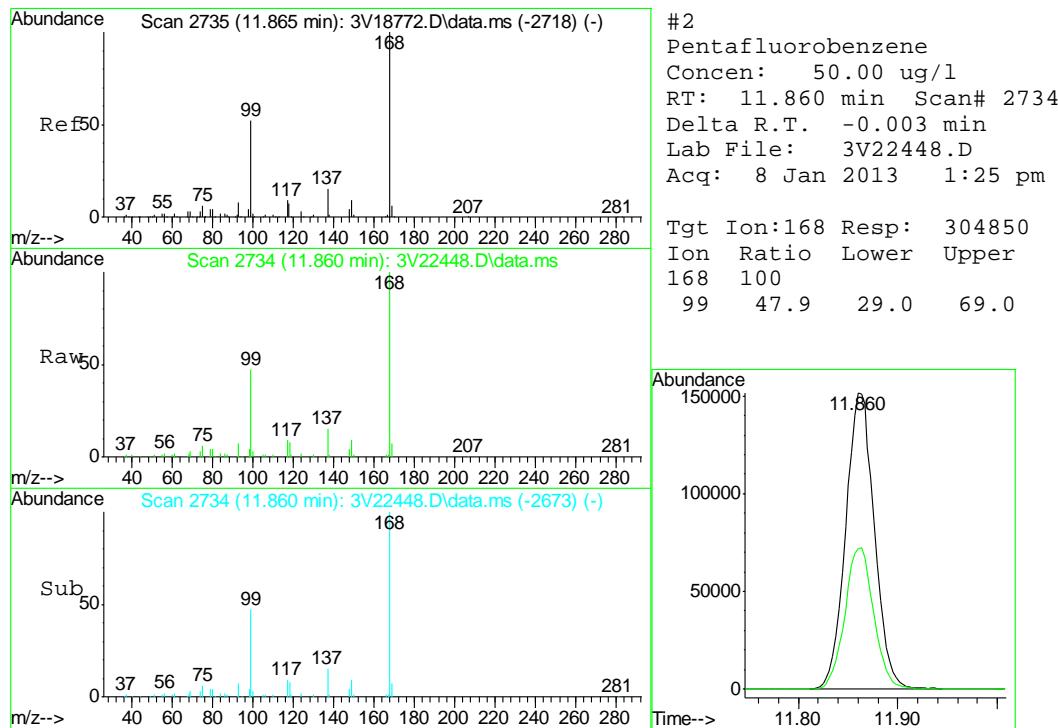
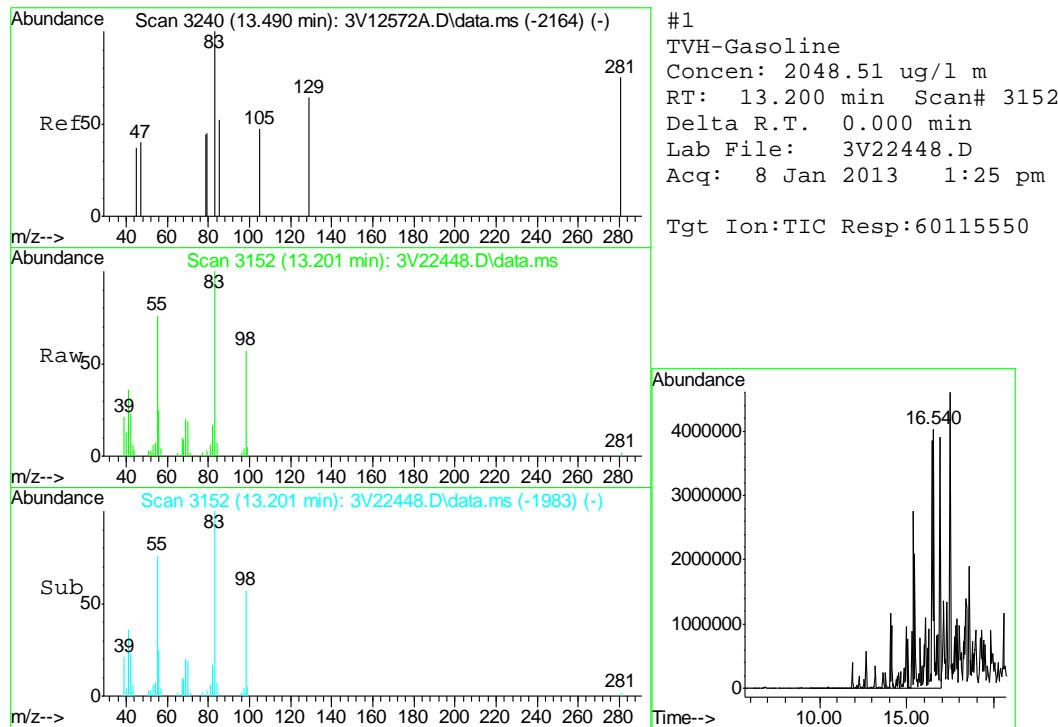
(#) = qualifier out of range (m) = manual integration (+) = signals summed

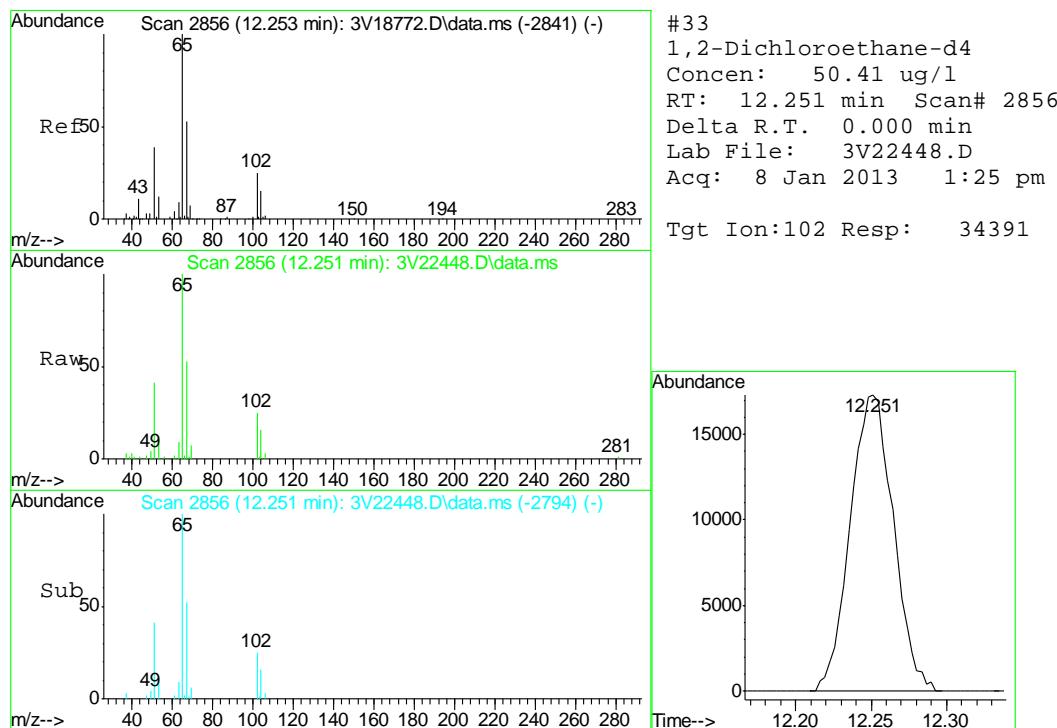
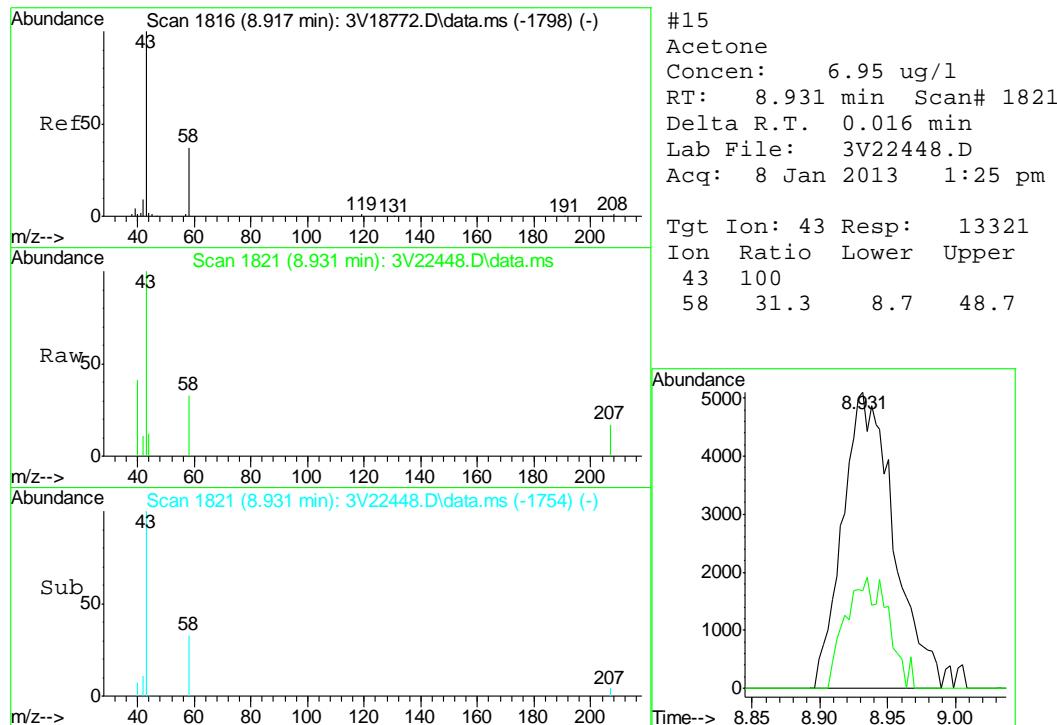
## Quantitation Report (QT Reviewed)

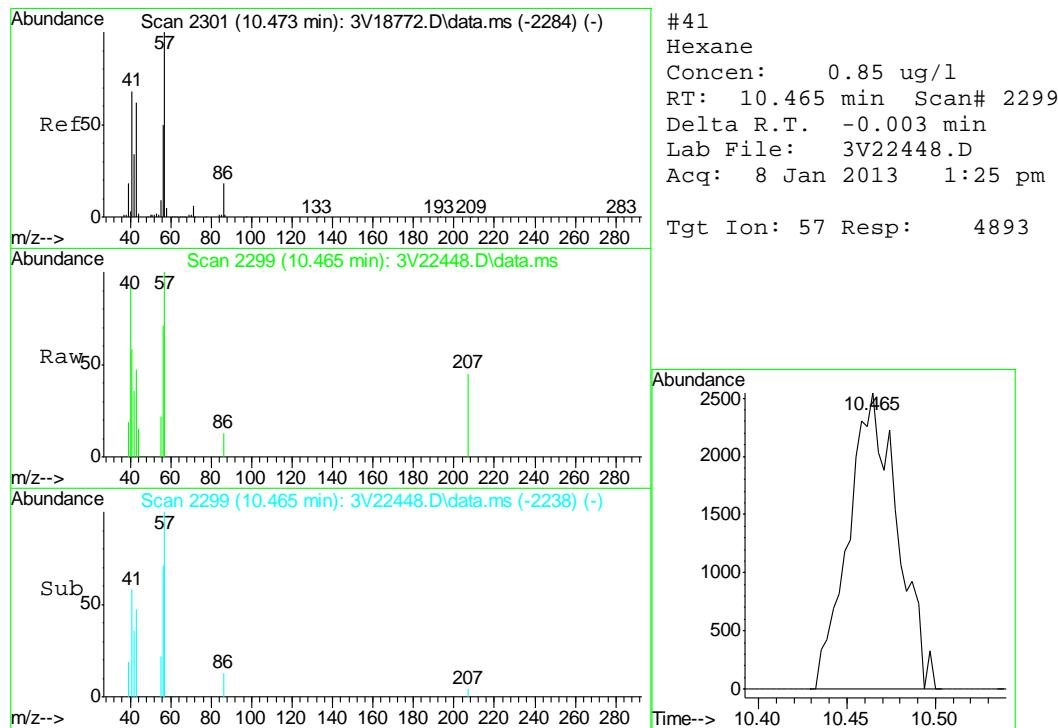
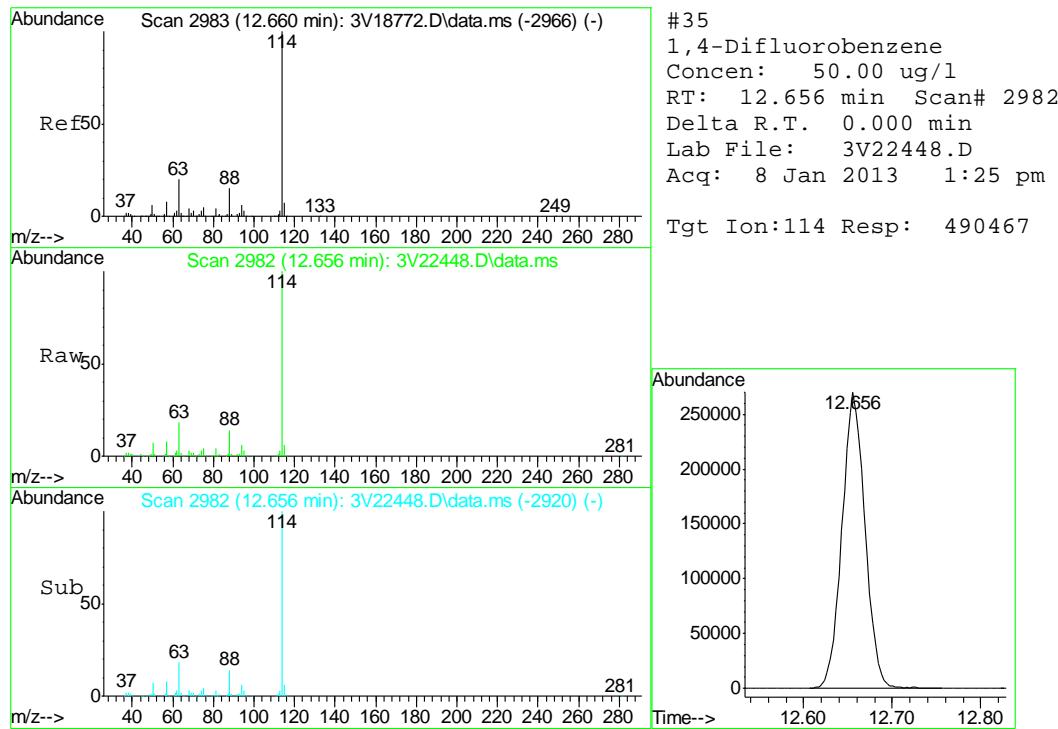
Data Path : C:\msdchem\1\DATA\V3010813.S\  
 Data File : 3V22448.D  
 Acq On : 8 Jan 2013 1:25 pm  
 Operator : BRETD  
 Sample : D42316-1  
 Misc : MS5203,V3V1321,5.022,,100,5,1  
 ALS Vial : 8 Sample Multiplier: 1

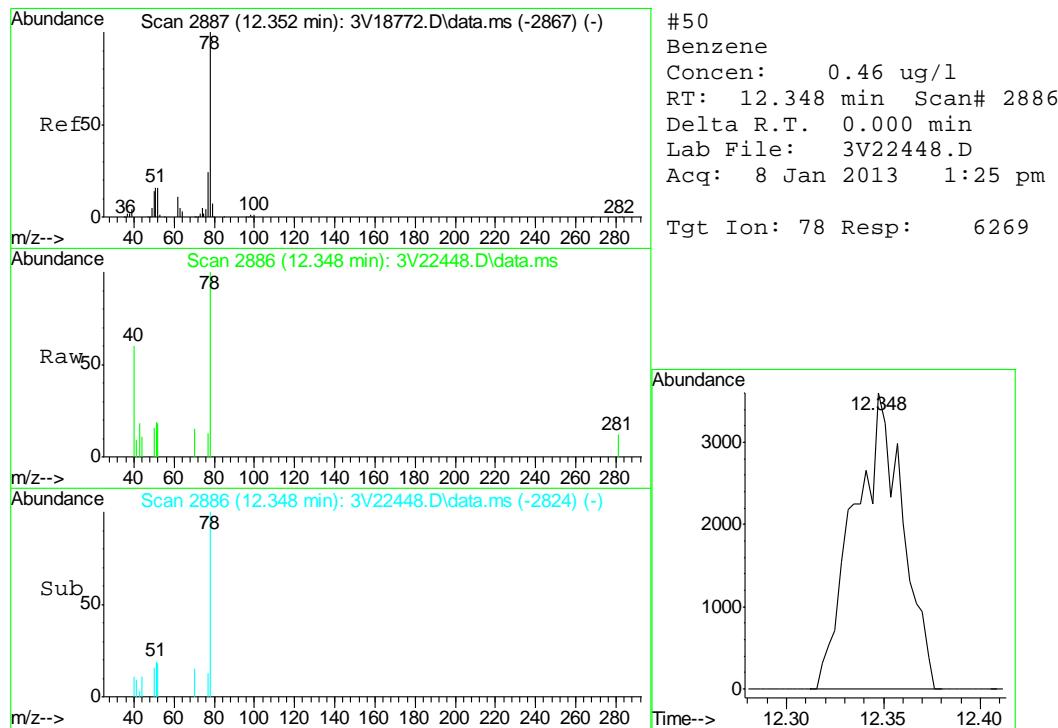
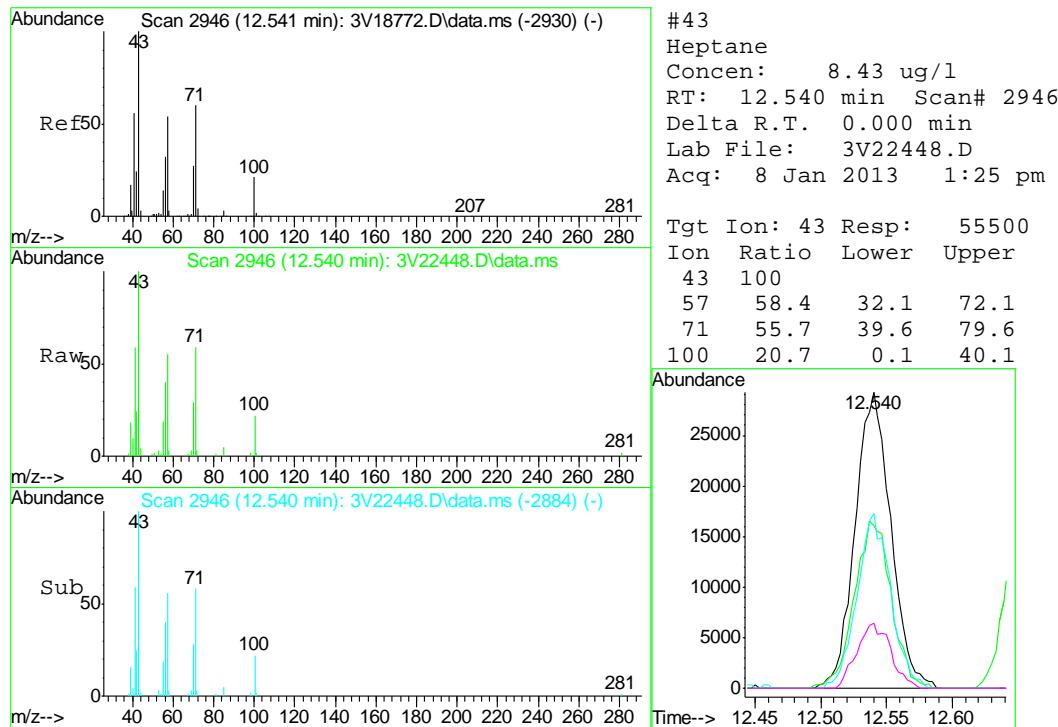
Quant Time: Jan 09 09:15:40 2013  
 Quant Method : C:\msdchem\1\METHODS\V3AP1299TVH1299SOIL.M  
 Quant Title : 8260  
 QLast Update : Thu Jan 03 11:40:16 2013  
 Response via : Initial Calibration

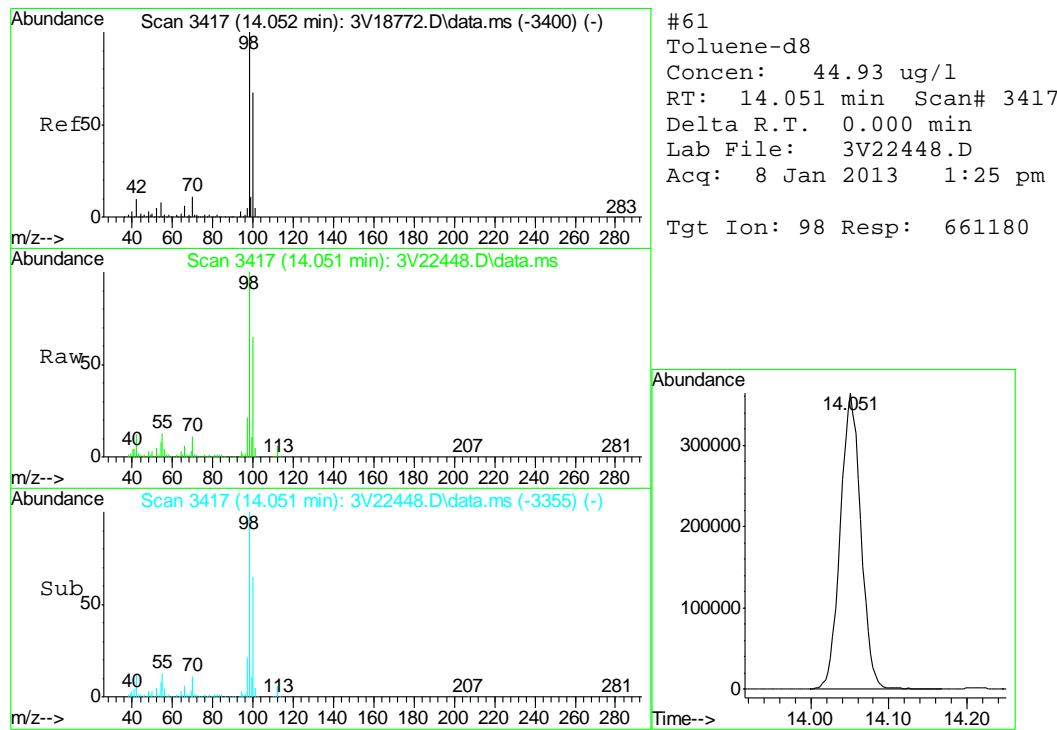
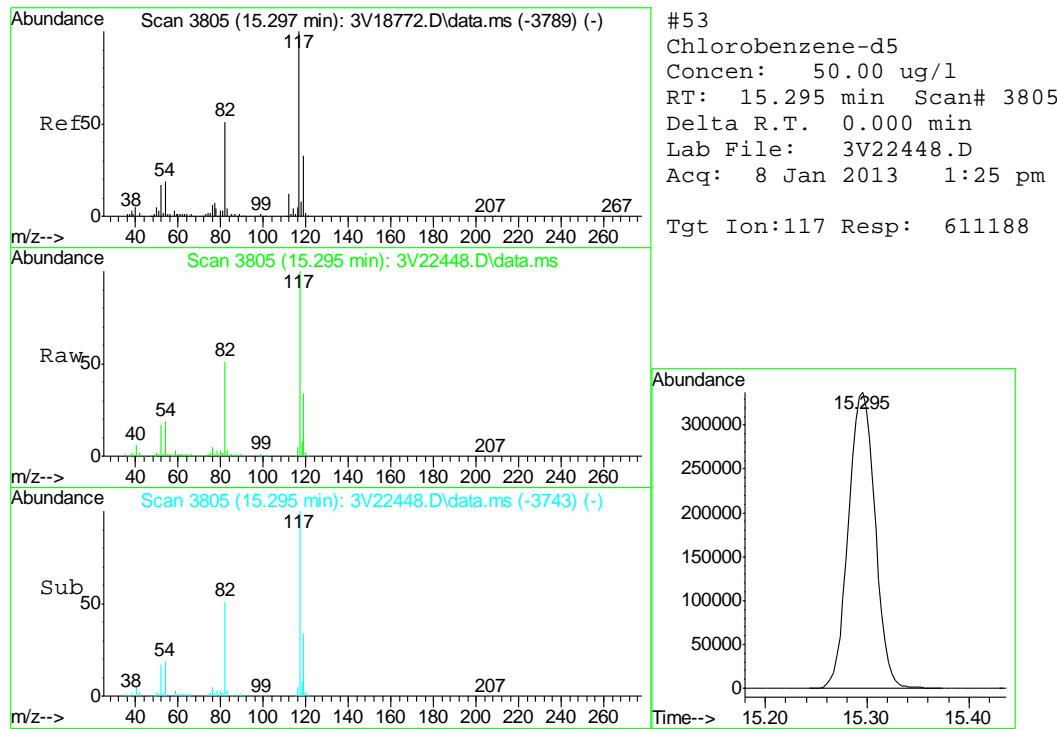


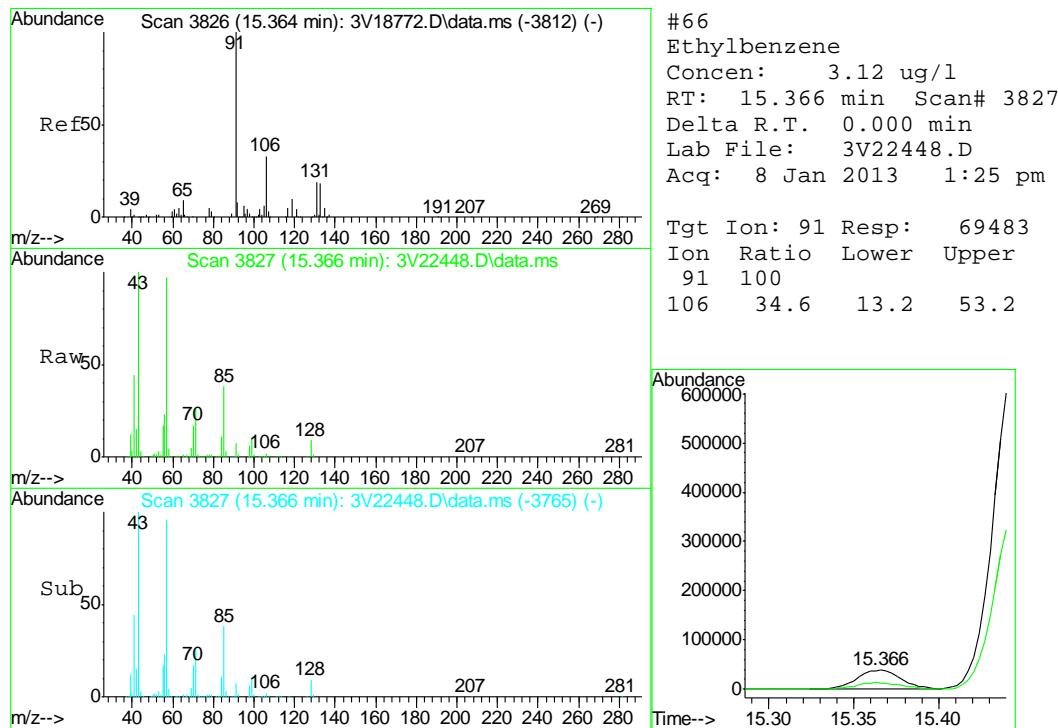
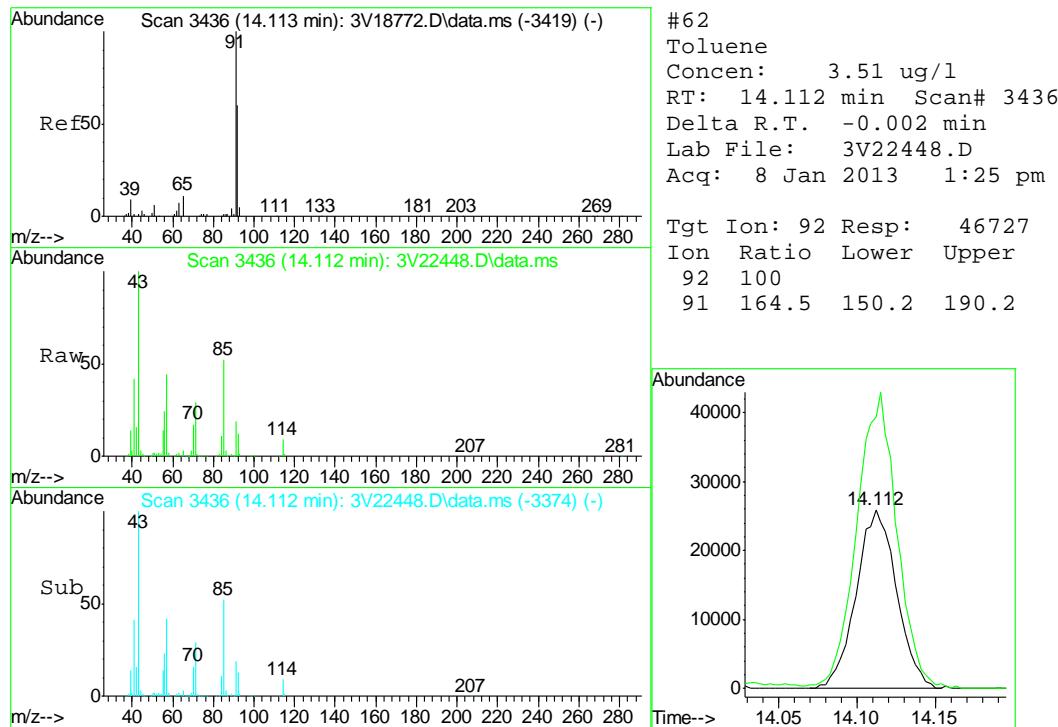


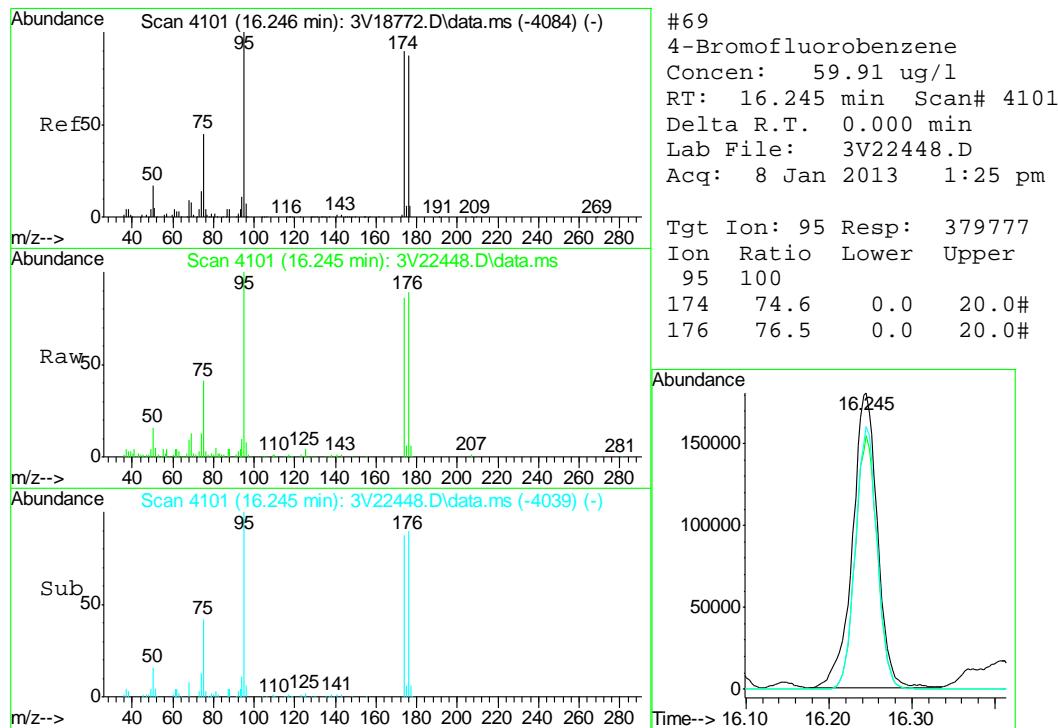
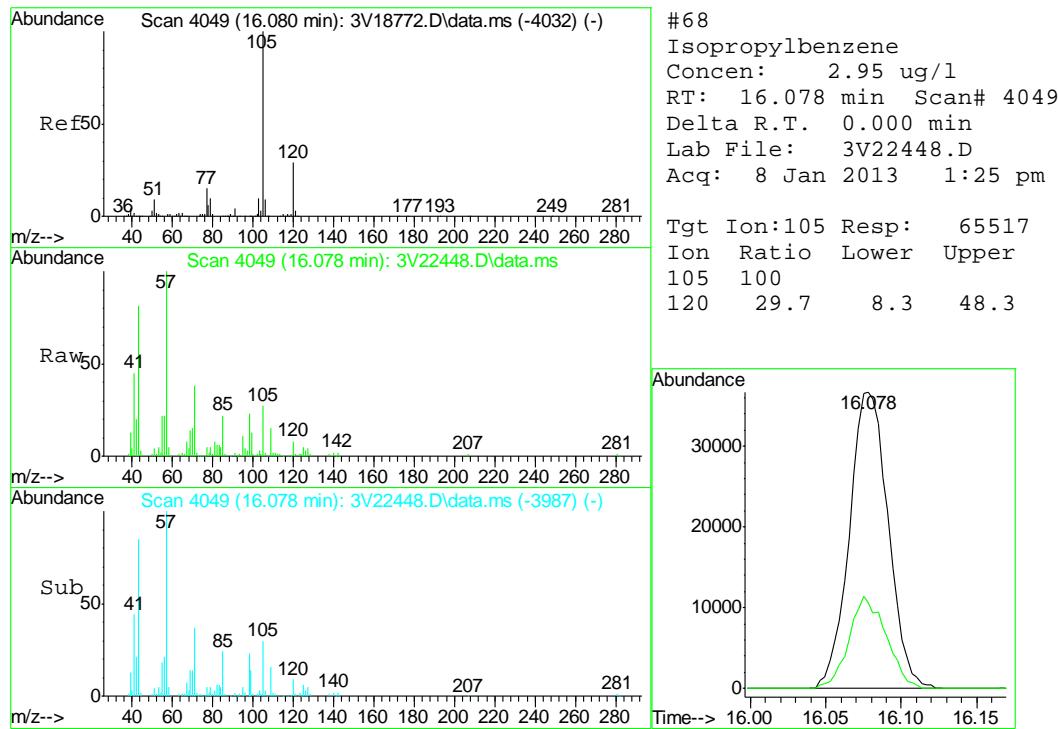


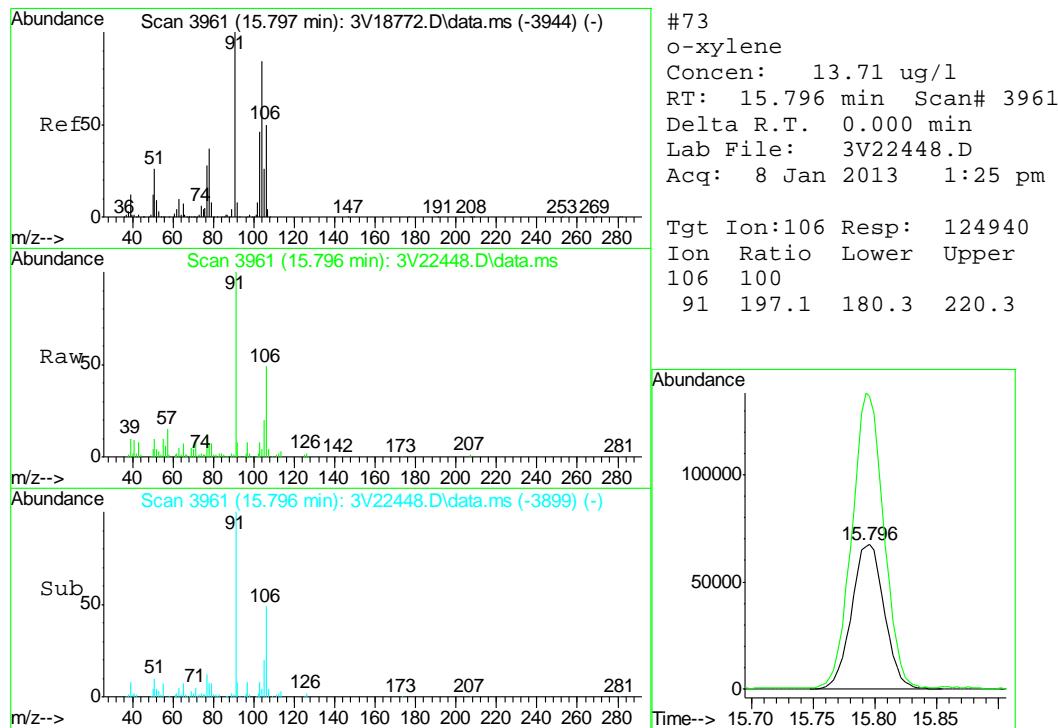
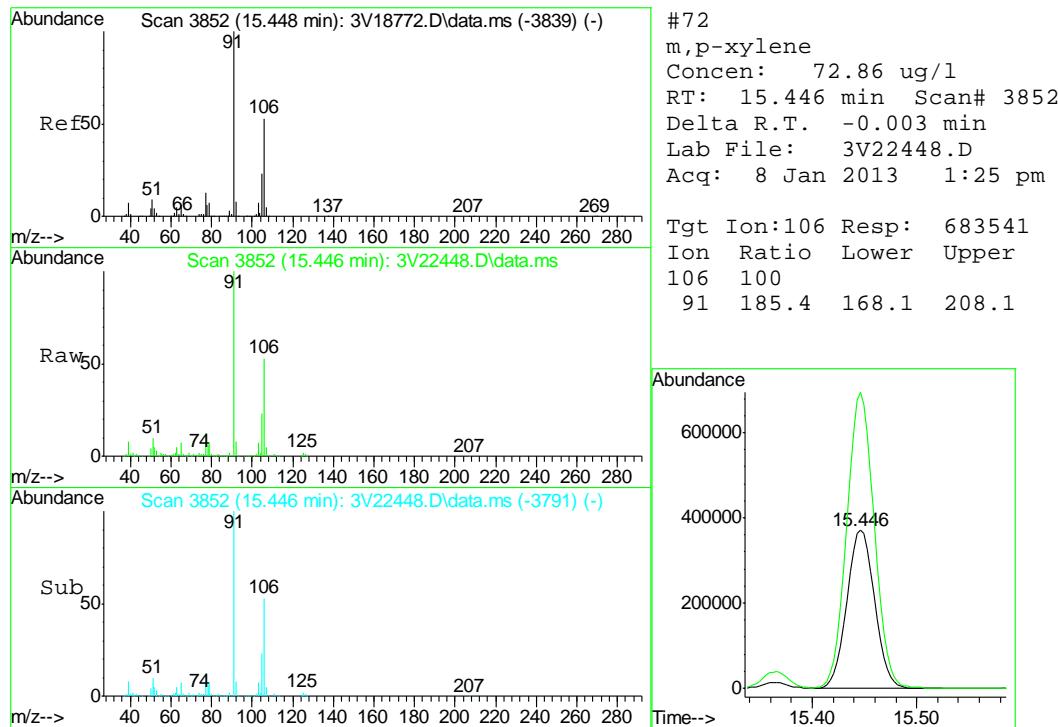


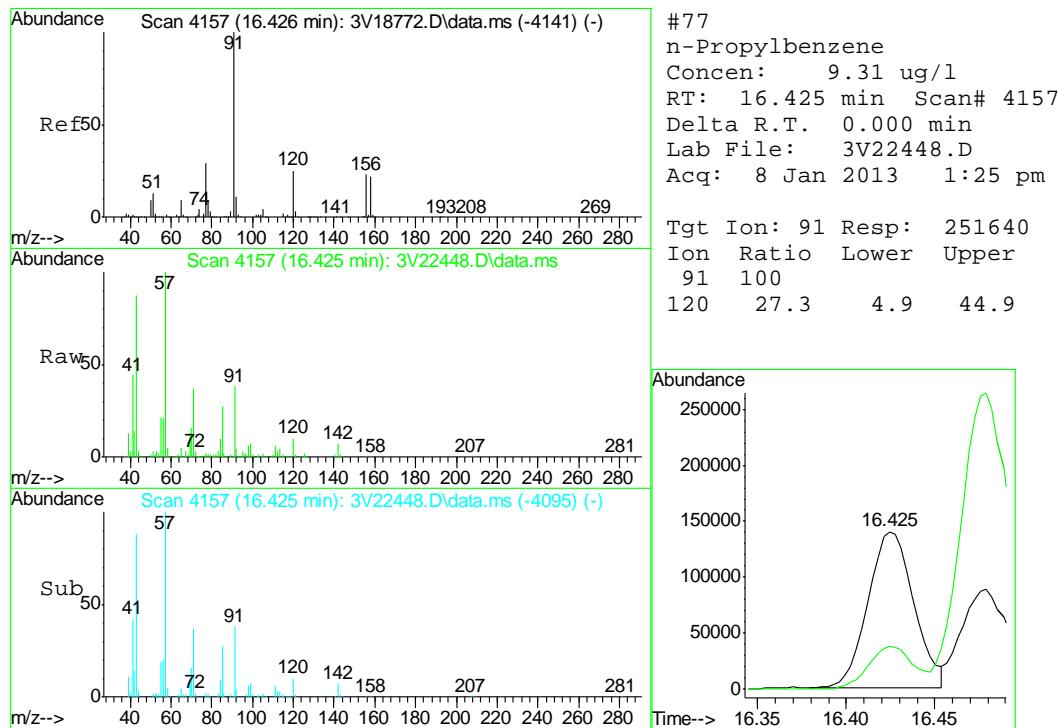
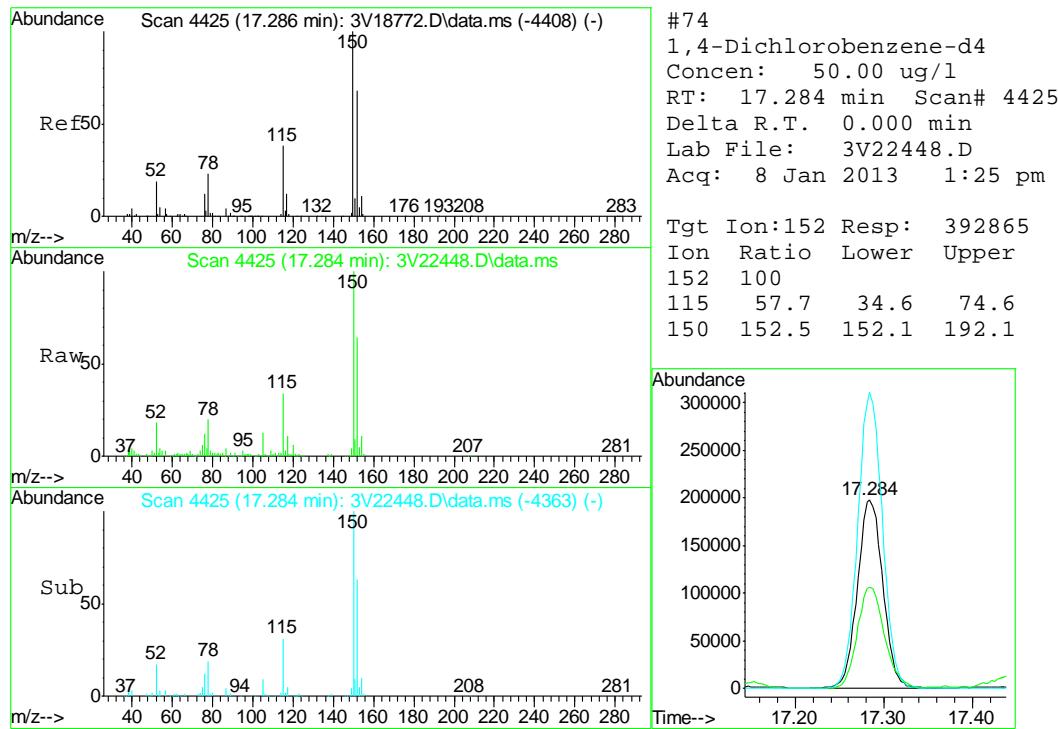


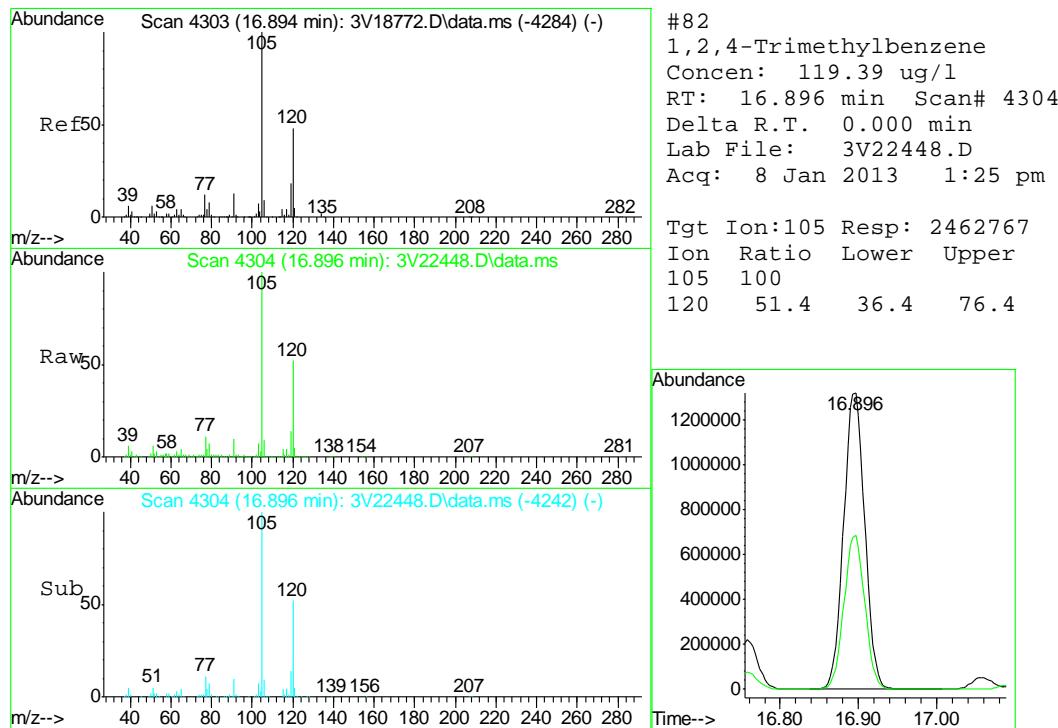
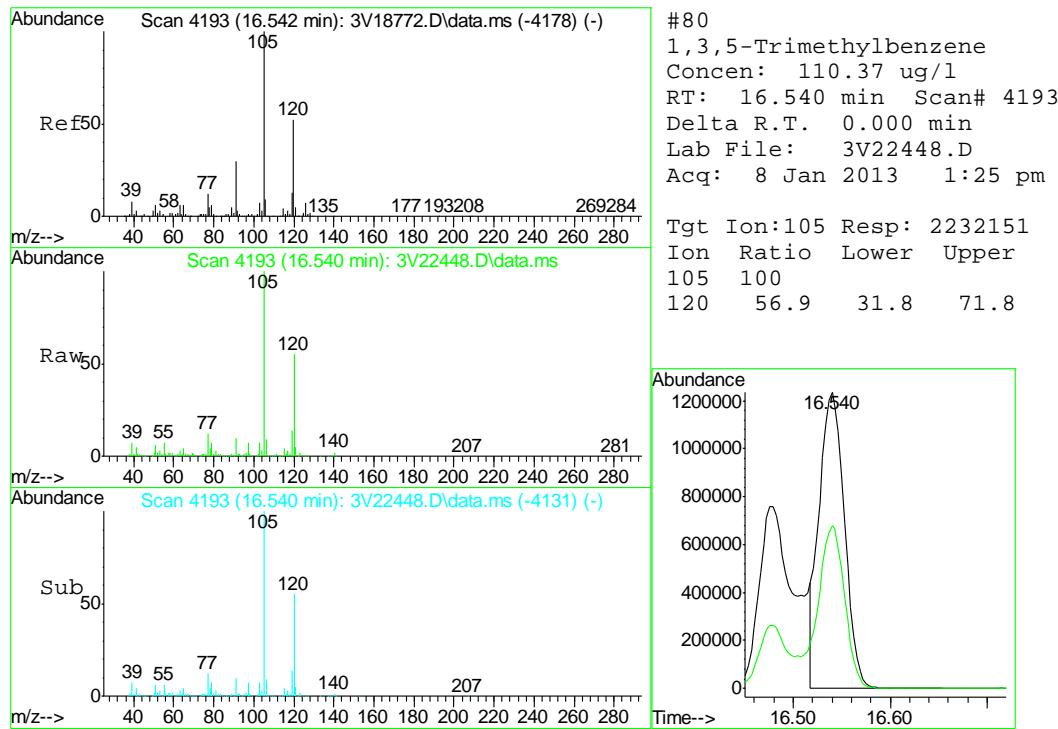


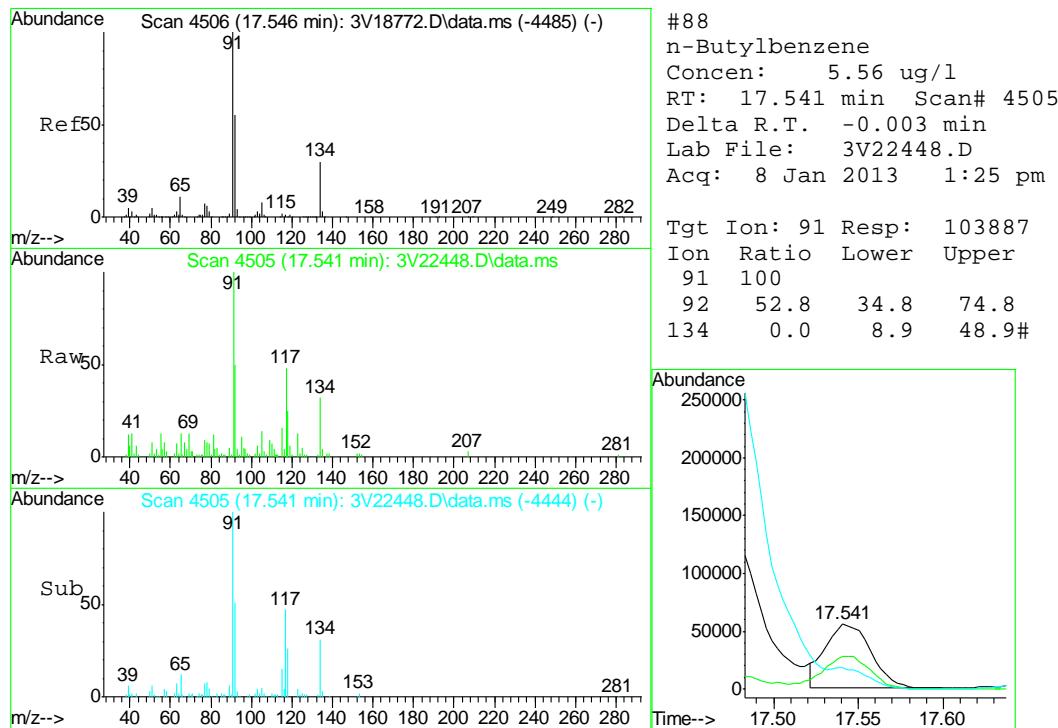
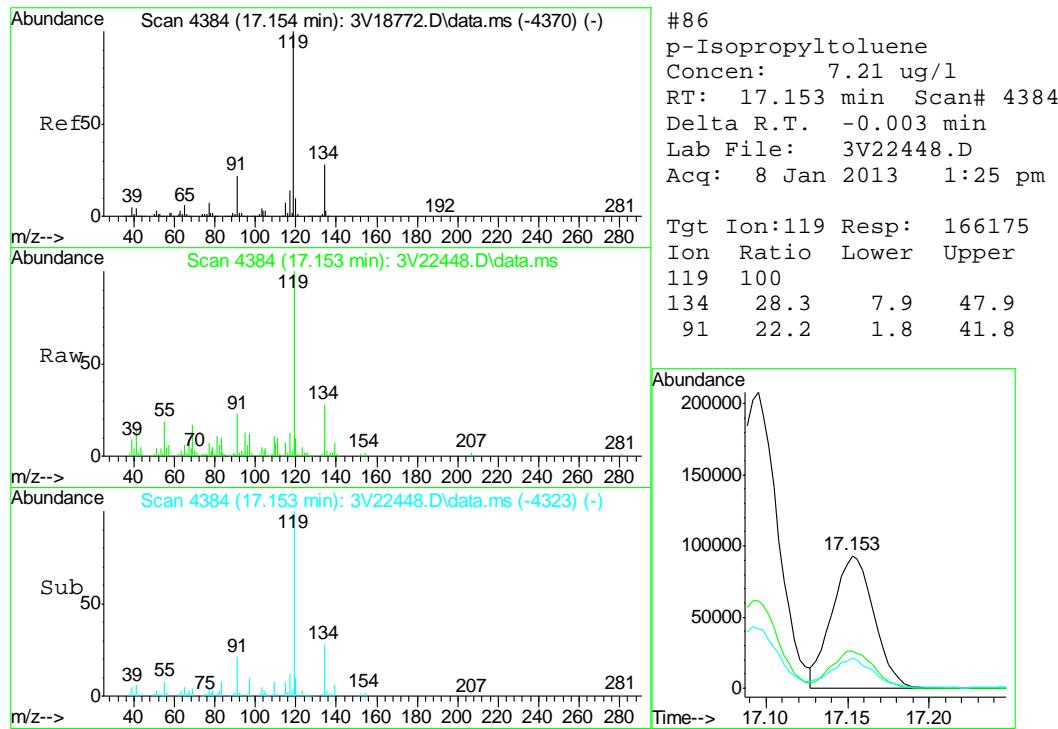


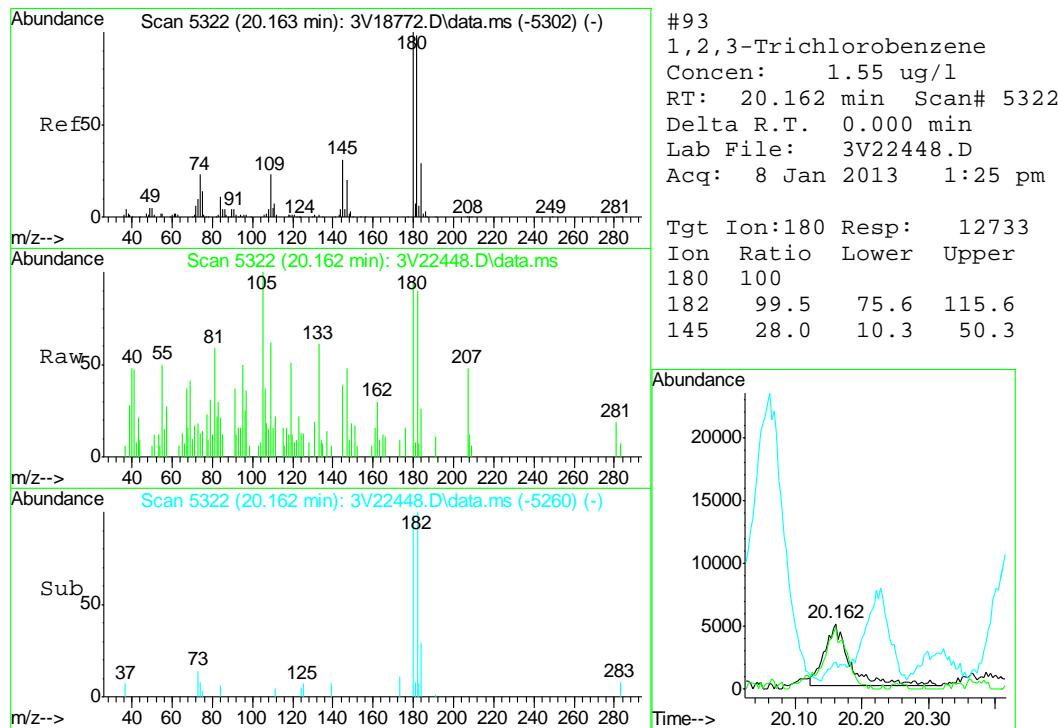
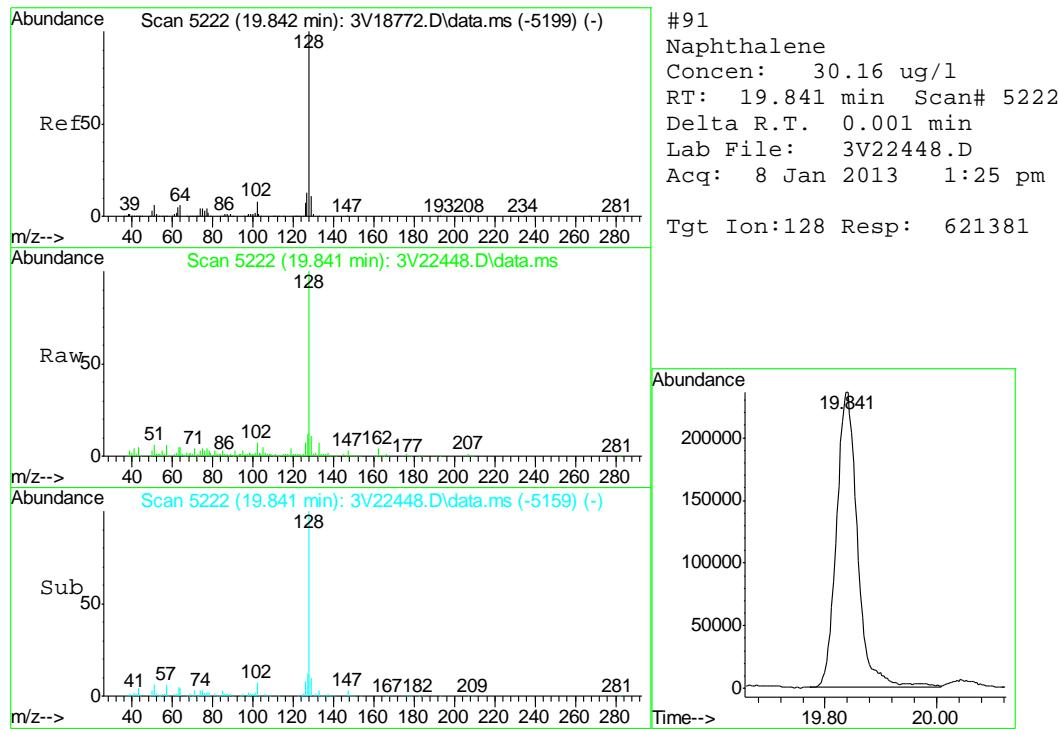


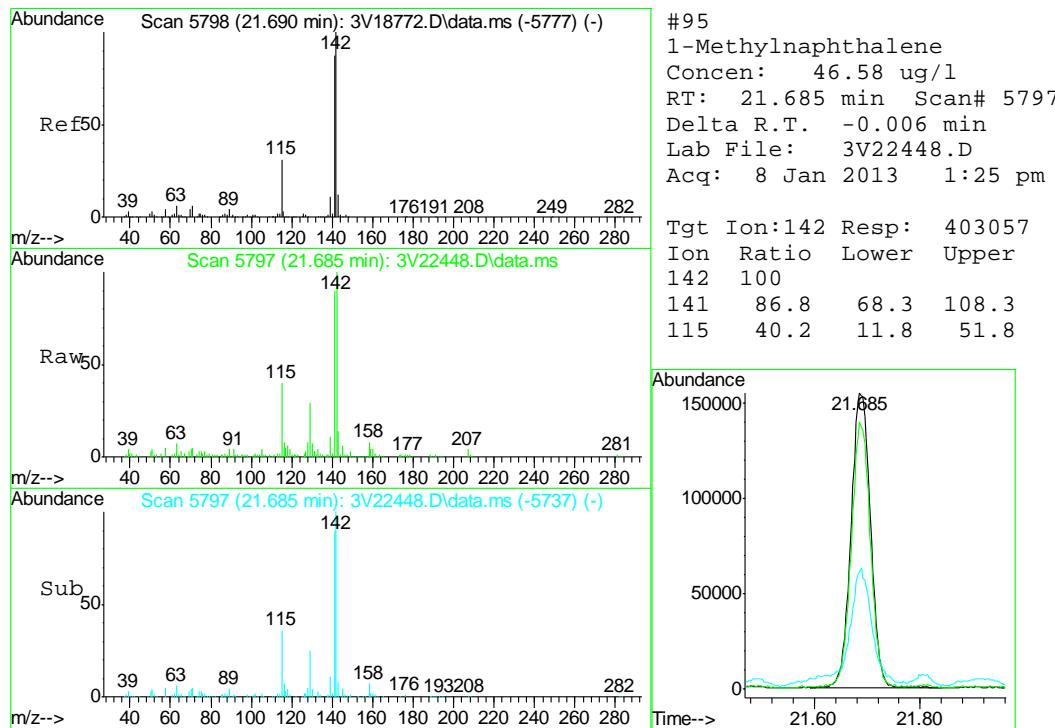
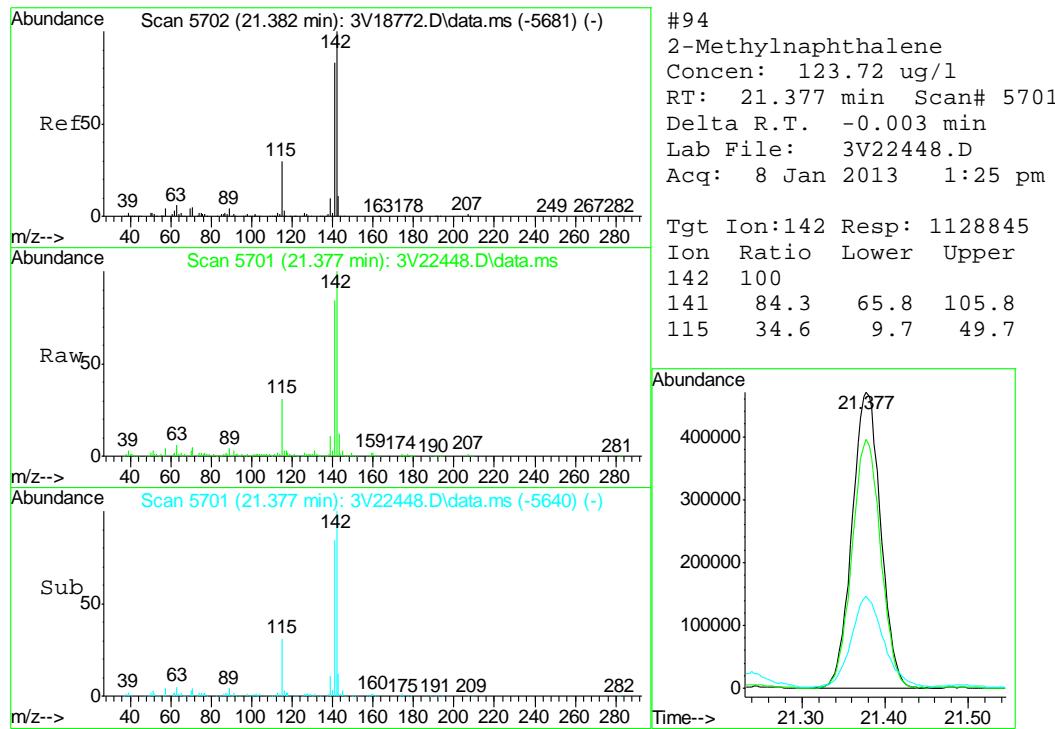












## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\V3010813.S\  
 Data File : 3V22443.D  
 Acq On : 8 Jan 2013 10:48 am  
 Operator : BRETD  
 Sample : MB  
 Misc : MS5203,V3V1321,5.00,,100,5,1  
 ALS Vial : 3 Sample Multiplier: 1

Quant Time: Jan 09 08:42:31 2013  
 Quant Method : C:\msdchem\1\METHODS\V3AP1299TVH1299SOIL.M  
 Quant Title : 8260  
 QLast Update : Thu Jan 03 11:40:16 2013  
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
2) Pentafluorobenzene	11.863	168	268628	50.00	ug/l	0.00
35) 1,4-Difluorobenzene	12.656	114	434746	50.00	ug/l	0.00
53) Chlorobenzene-d5	15.296	117	540657	50.00	ug/l	0.00
74) 1,4-Dichlorobenzene-d4	17.285	152	324831	50.00	ug/l	0.00

System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	12.255	102	32637	54.29	ug/l	0.00
Spiked Amount	50.000	Range	70 - 130	Recovery	=	108.58%
61) Toluene-d8	14.051	98	556286	42.74	ug/l	0.00
Spiked Amount	50.000	Range	70 - 130	Recovery	=	85.48%
69) 4-Bromofluorobenzene	16.242	95	267407	47.69	ug/l	0.00
Spiked Amount	50.000	Range	70 - 130	Recovery	=	95.38%

Target Compounds					Qvalue
1) TVH-Gasoline	13.200	TIC	87958m	3.00	ug/l
15) Acetone	8.932	43	9364	4.35	ug/l
70) Bromobenzene	16.431	156	1458	0.25	ug/l
71) Styrene	15.793	104	2183	0.57	ug/l
79) 4-Chlorotoluene	16.646	91	2958	0.20	ug/l
84) 1,3-Dichlorobenzene	17.240	146	3389	0.30	ug/l
87) 1,2-Dichlorobenzene	17.692	146	3642	0.33	ug/l
88) n-Butylbenzene	17.551	91	3434	0.22	ug/l
90) 1,2,4-Trichlorobenzene	19.463	180	3957	0.57	ug/l
91) Naphthalene	19.841	128	19171	2.11	ug/l
93) 1,2,3-Trichlorobenzene	20.159	180	4031	0.59	ug/l
94) 2-Methylnaphthalene	21.371	142	14756	1.96	ug/l
95) 1-Methylnaphthalene	21.689	142	12687	1.77	ug/l

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

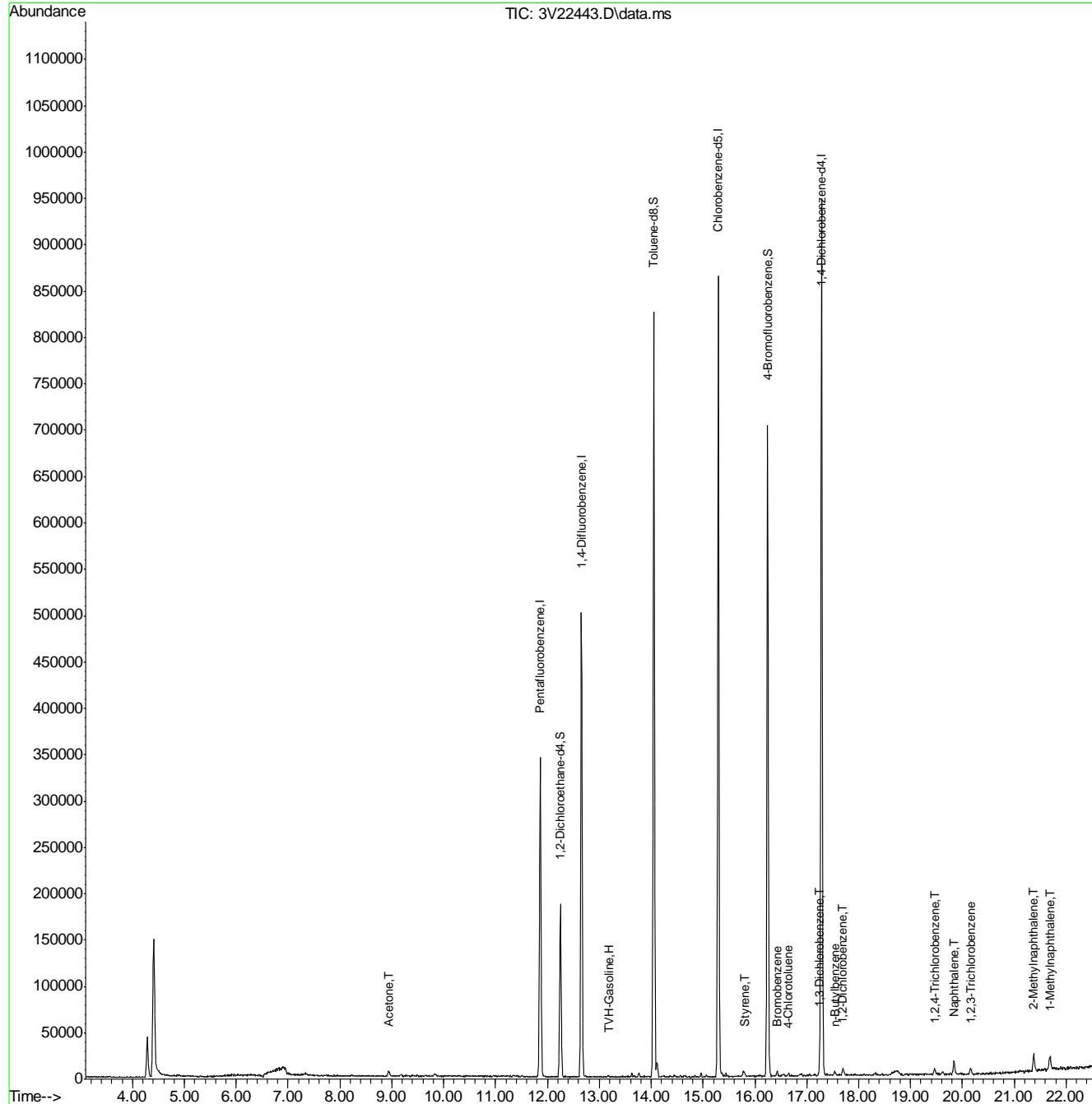
7.2.1

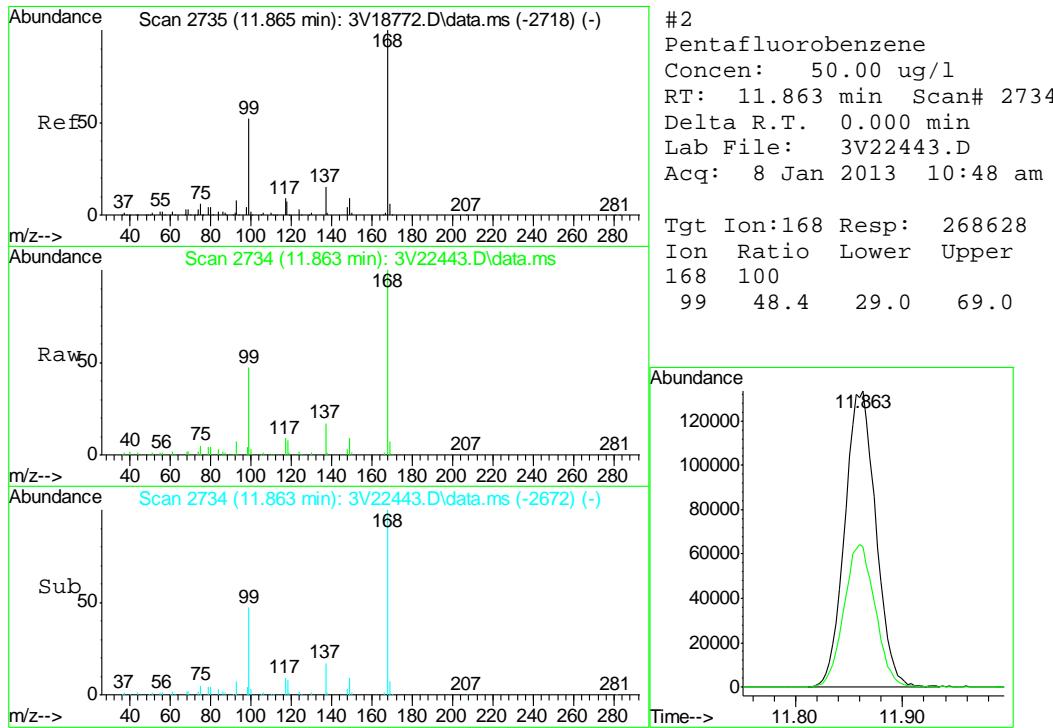
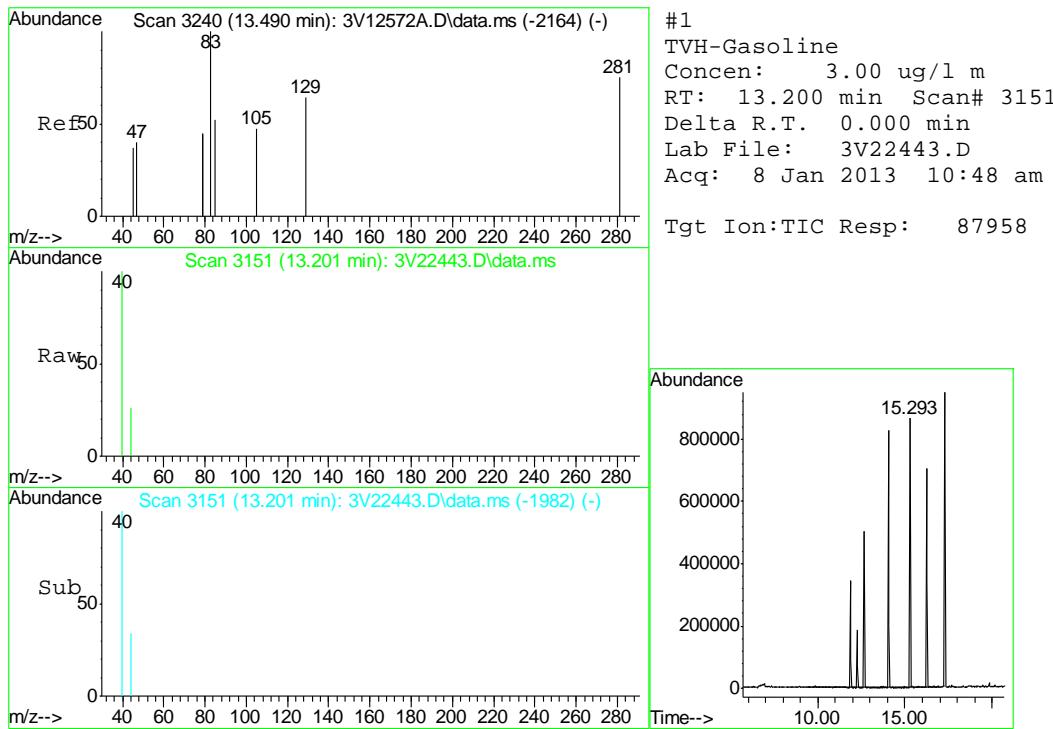
7

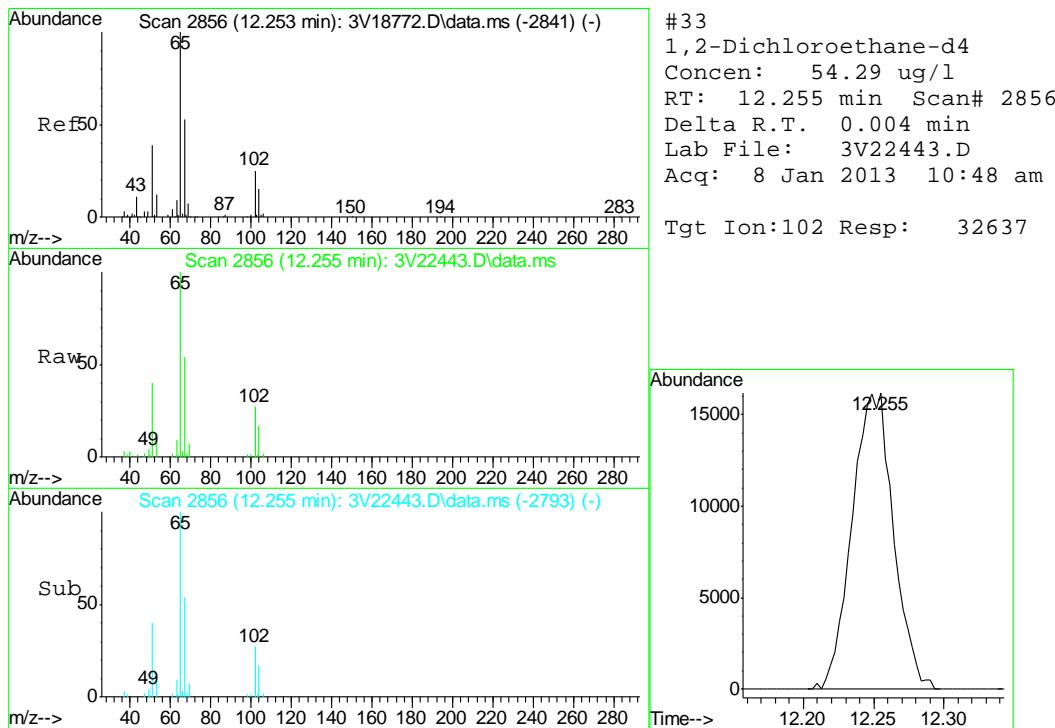
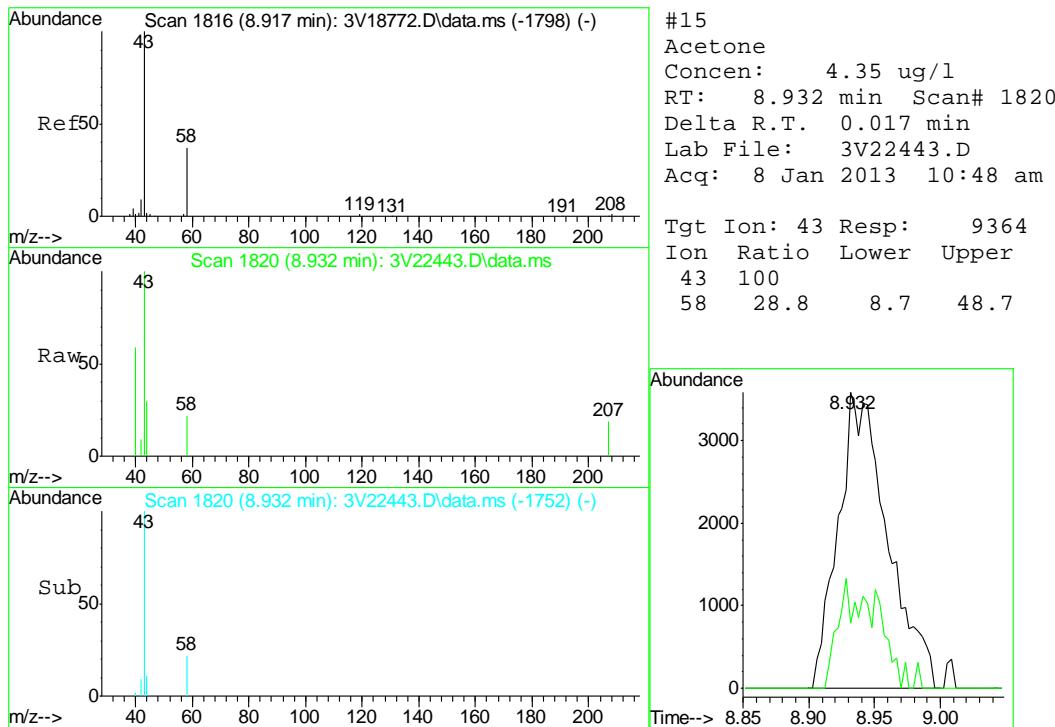
## Quantitation Report (QT Reviewed)

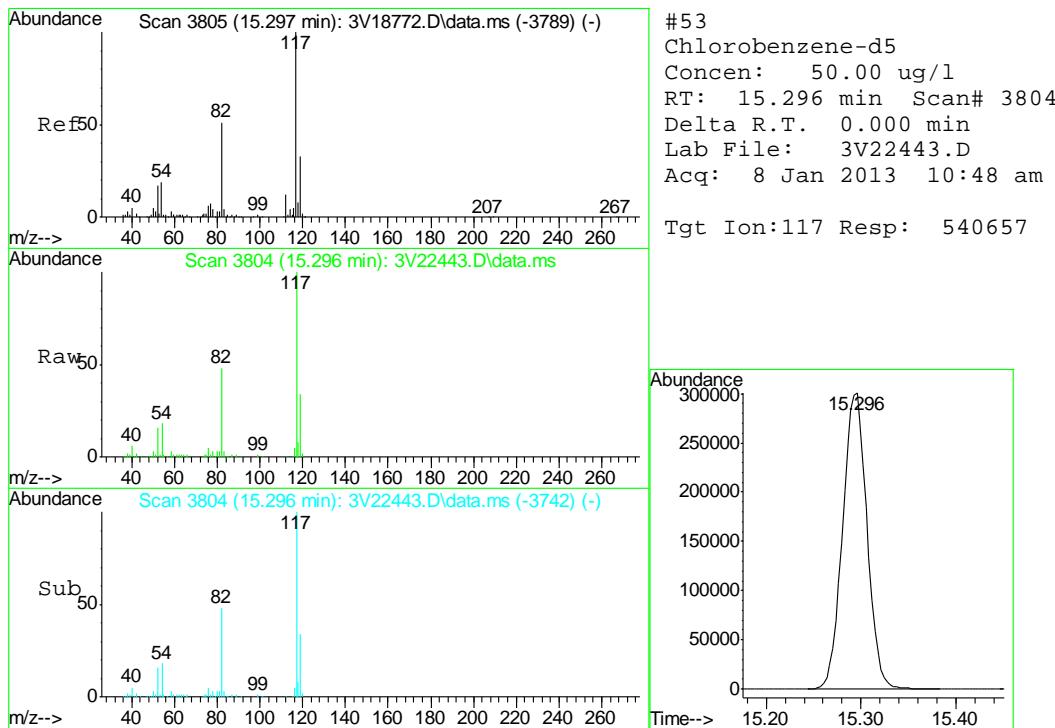
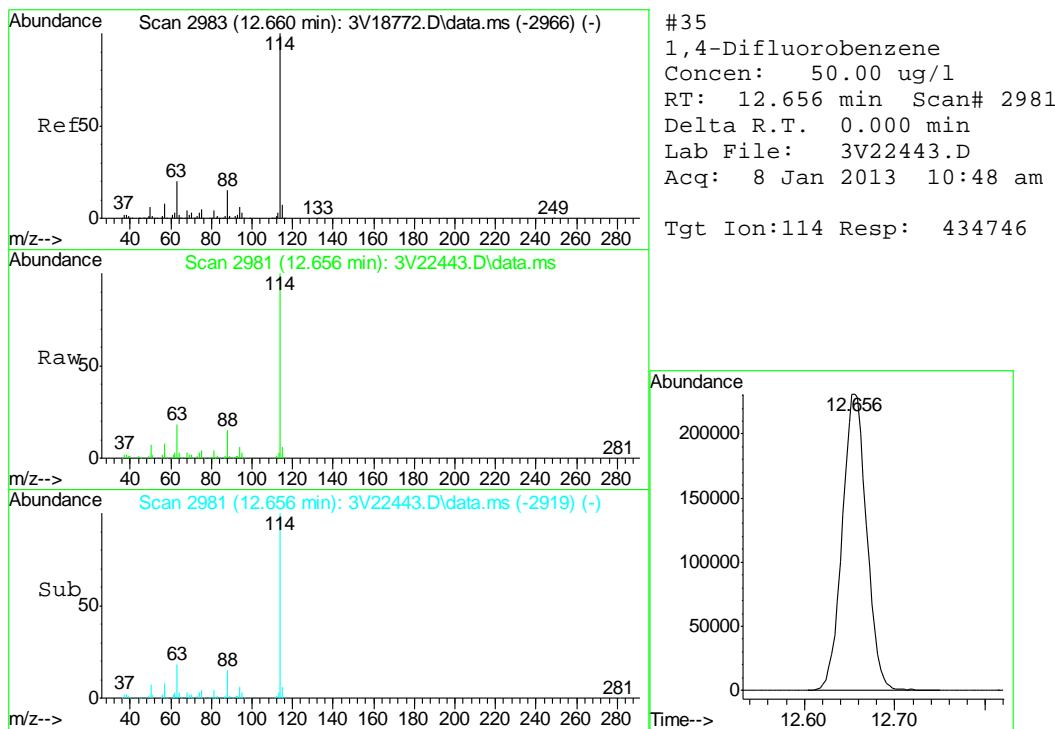
Data Path : C:\msdchem\1\DATA\V3010813.S\  
 Data File : 3V22443.D  
 Acq On : 8 Jan 2013 10:48 am  
 Operator : BRETD  
 Sample : MB  
 Misc : MS5203,V3V1321,5.00,,100,5,1  
 ALS Vial : 3 Sample Multiplier: 1

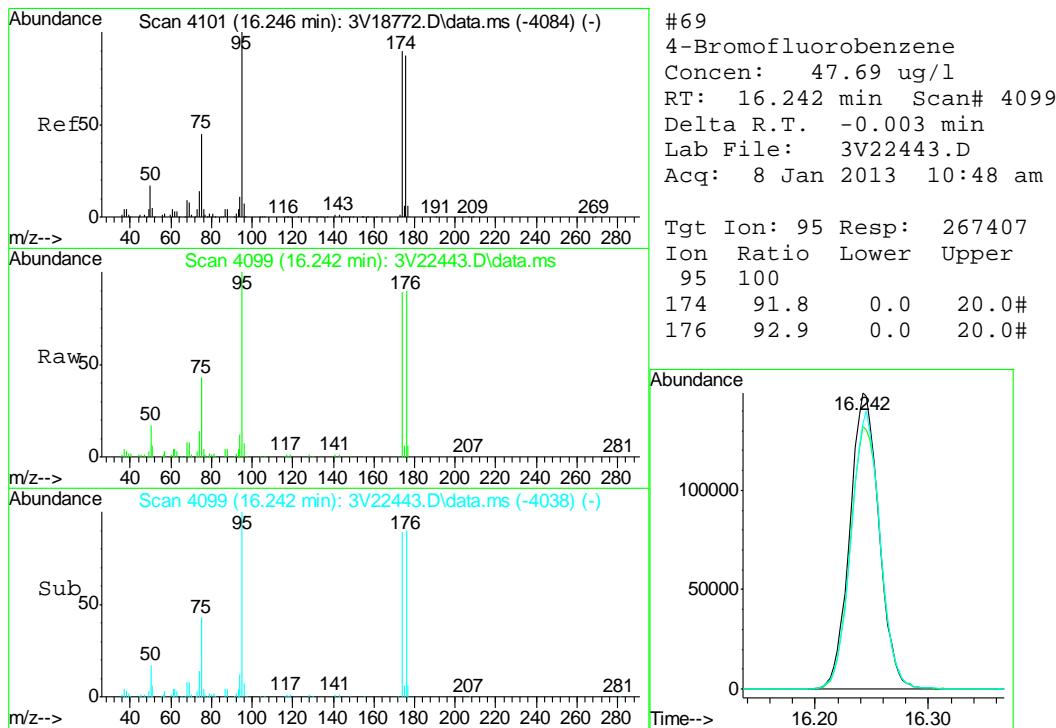
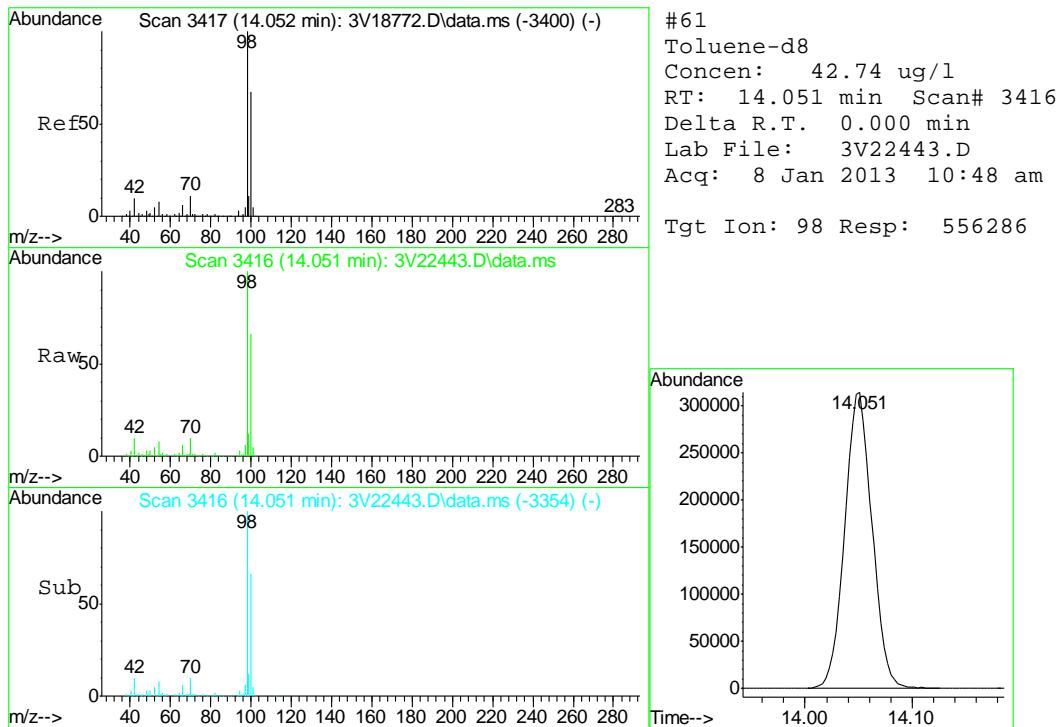
Quant Time: Jan 09 08:42:31 2013  
 Quant Method : C:\msdchem\1\METHODS\V3AP1299TVH1299SOIL.M  
 Quant Title : 8260  
 QLast Update : Thu Jan 03 11:40:16 2013  
 Response via : Initial Calibration

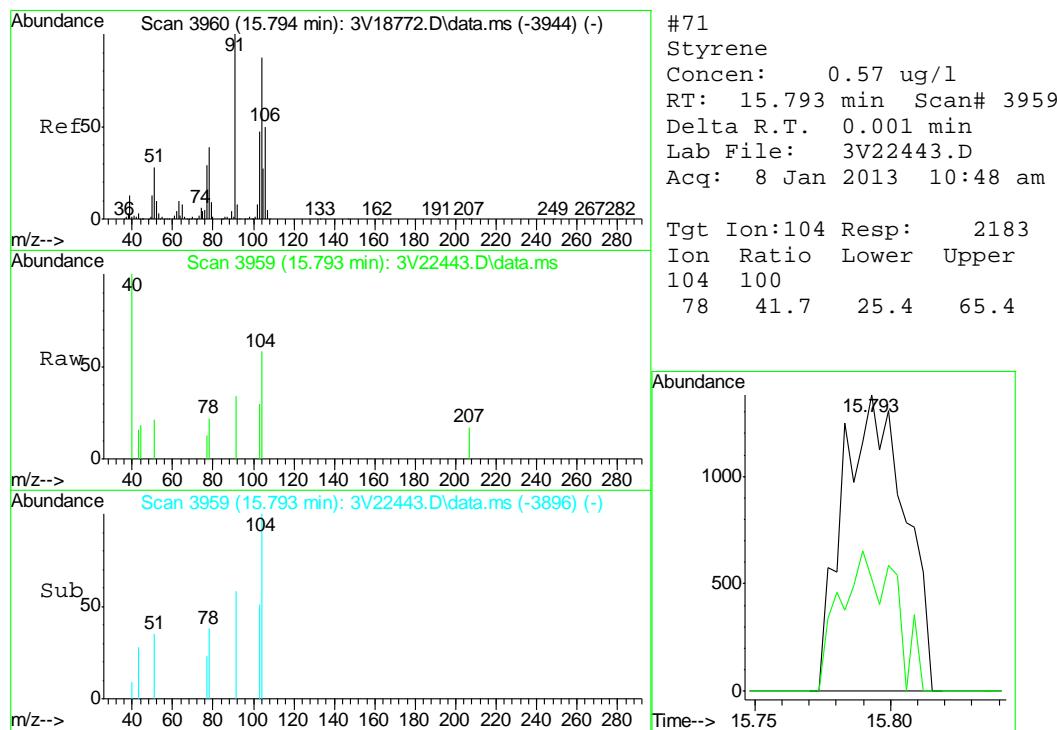
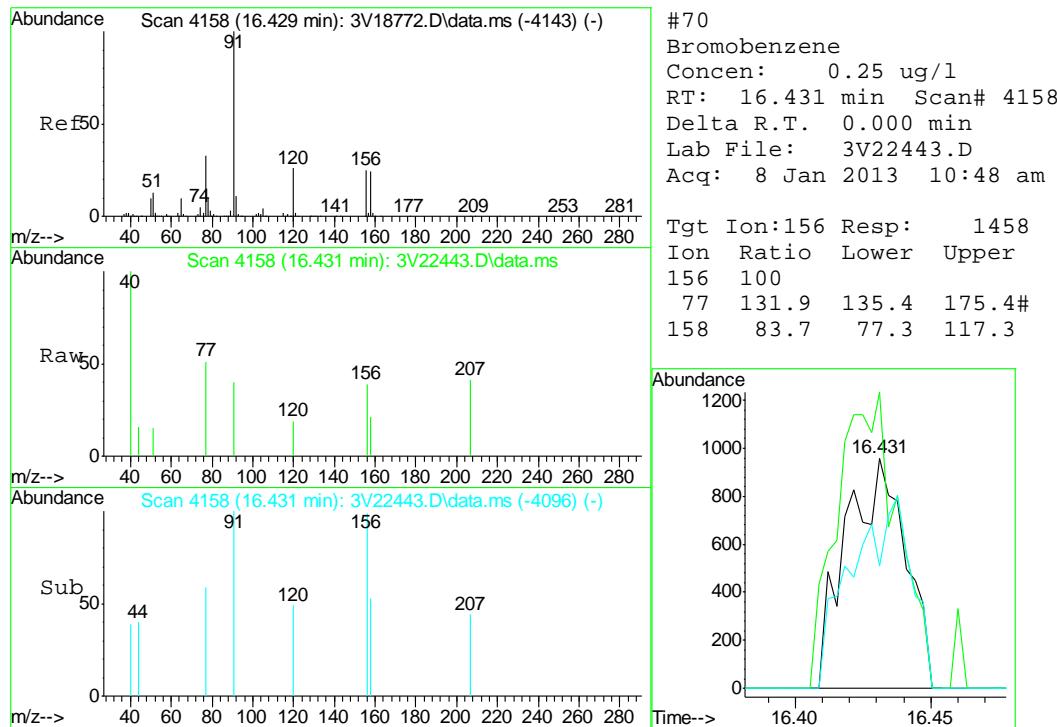


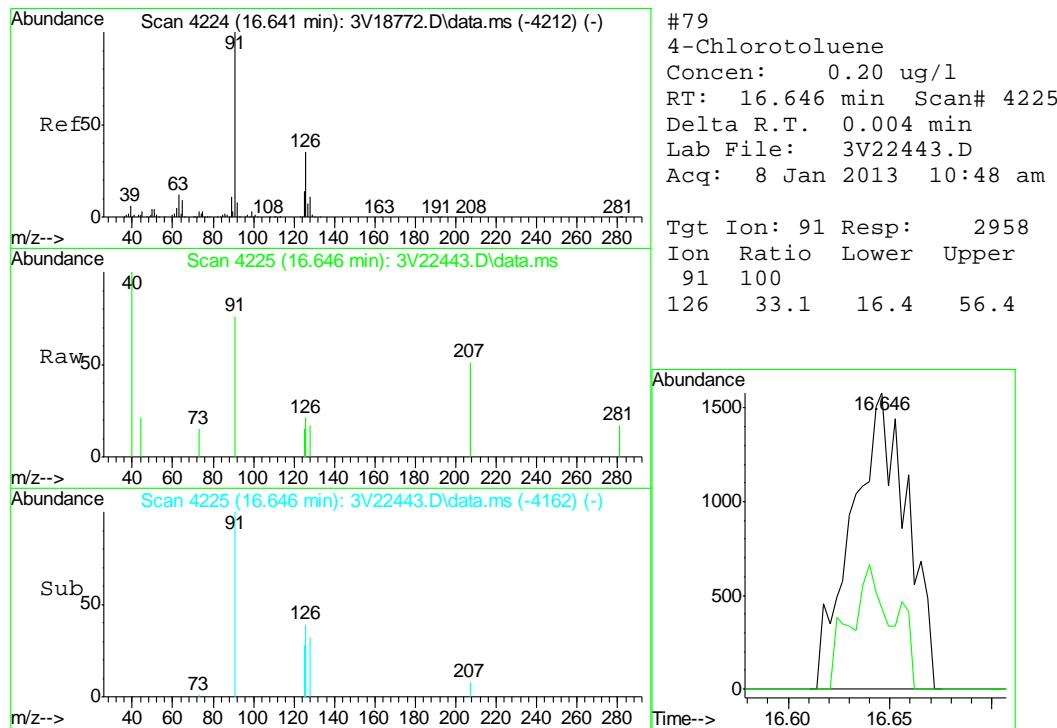
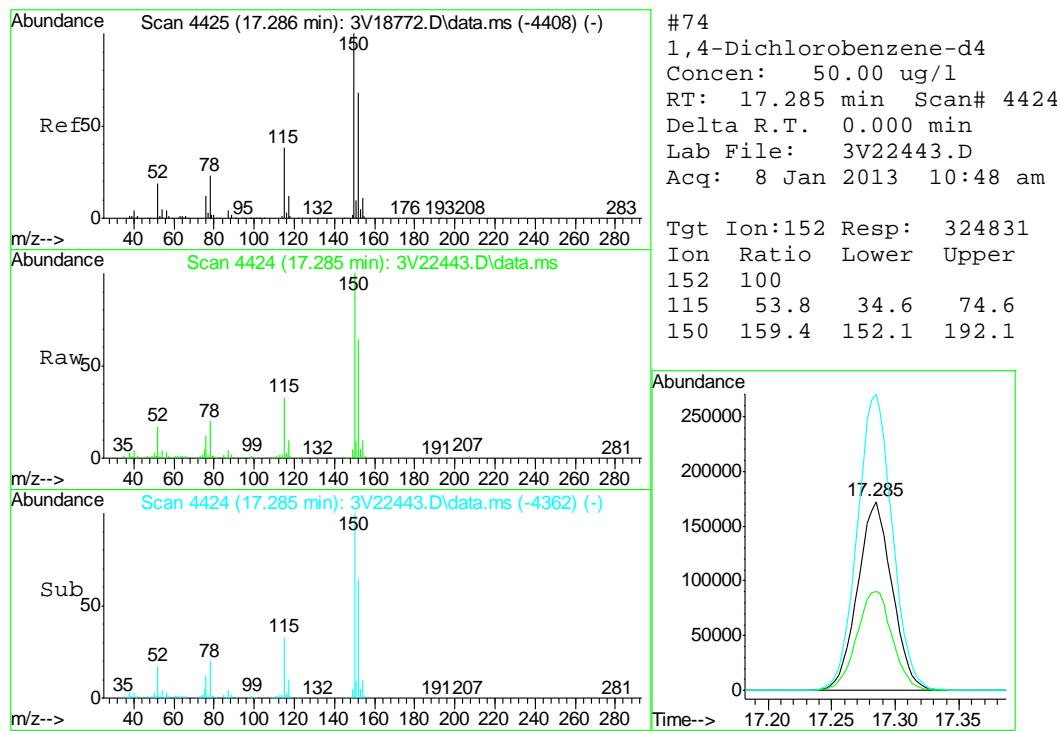


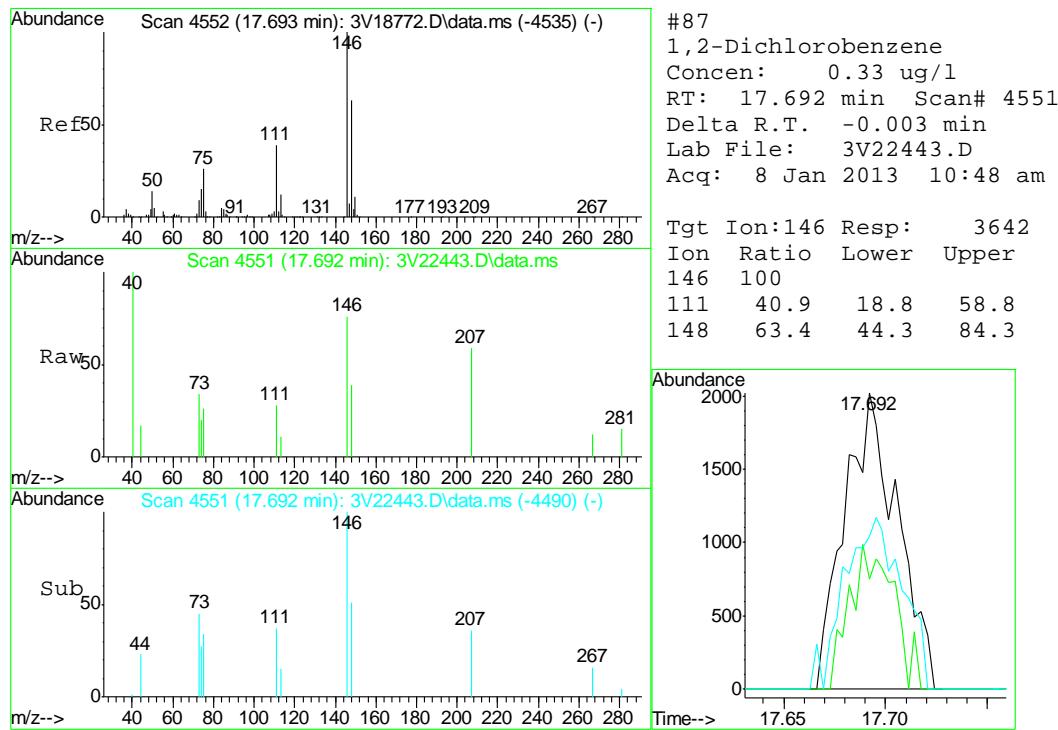
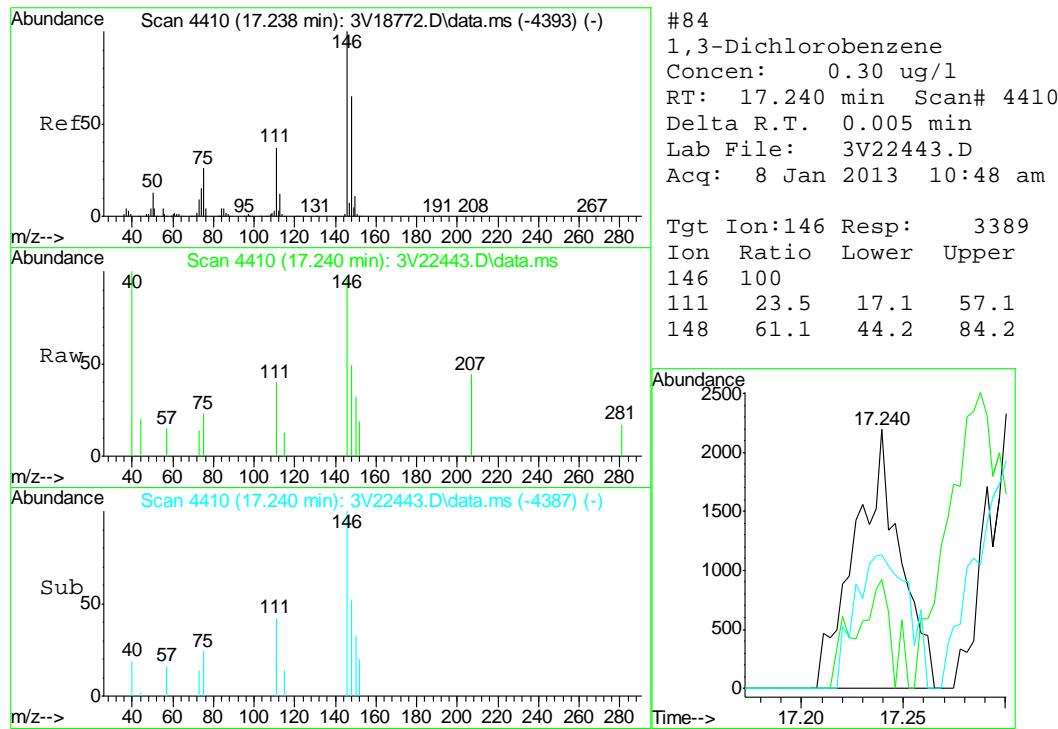


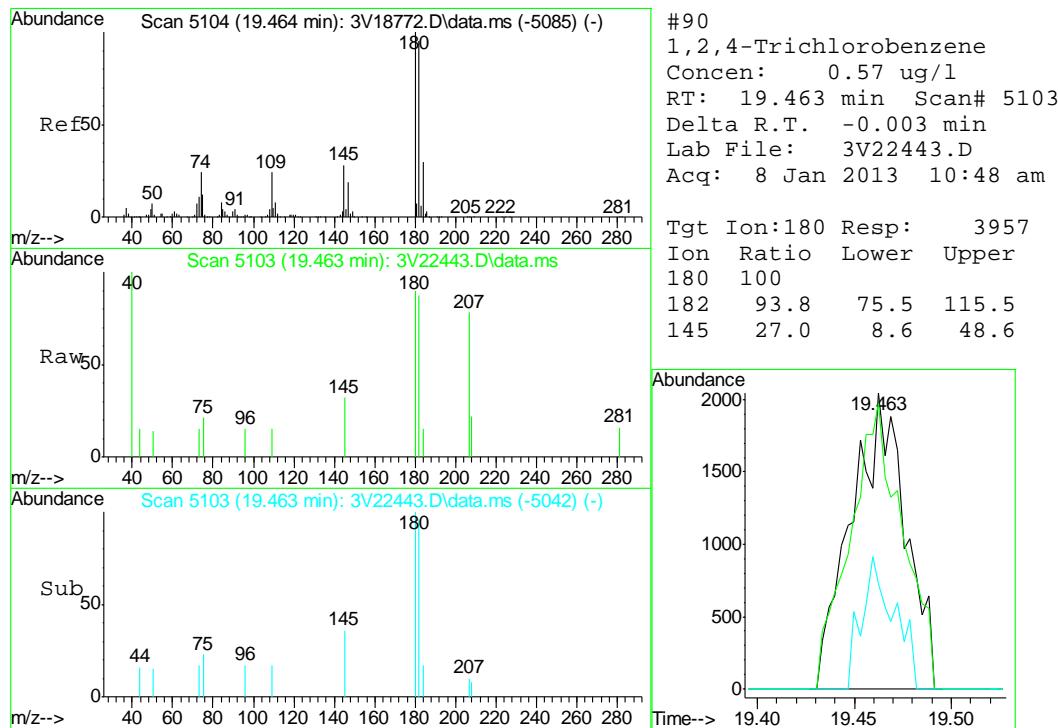
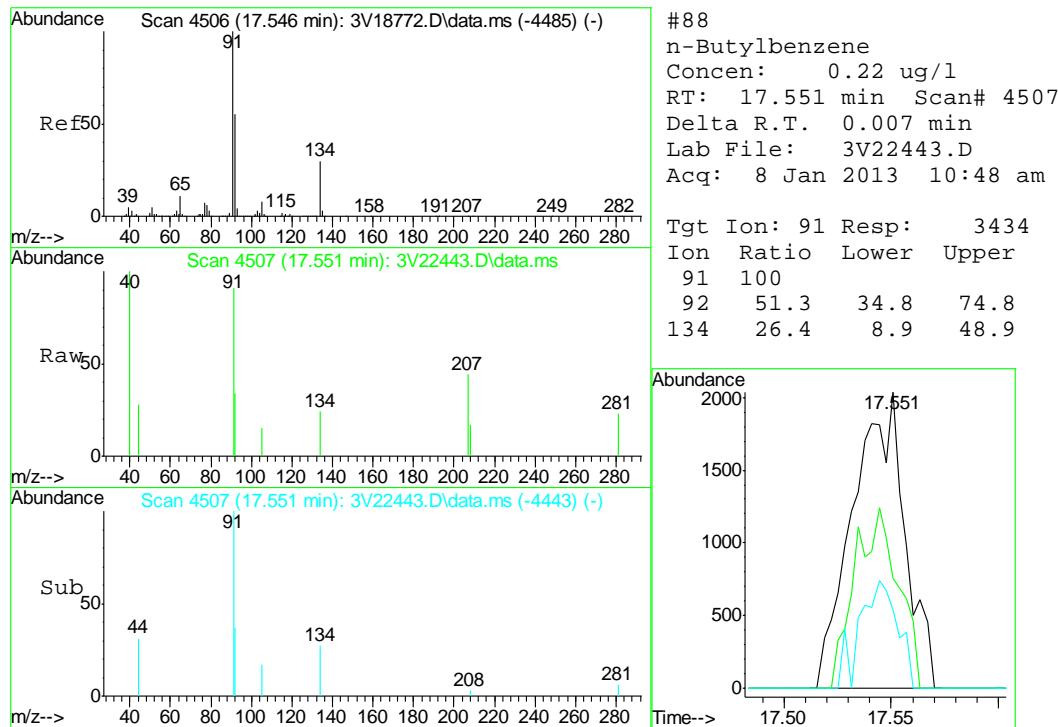


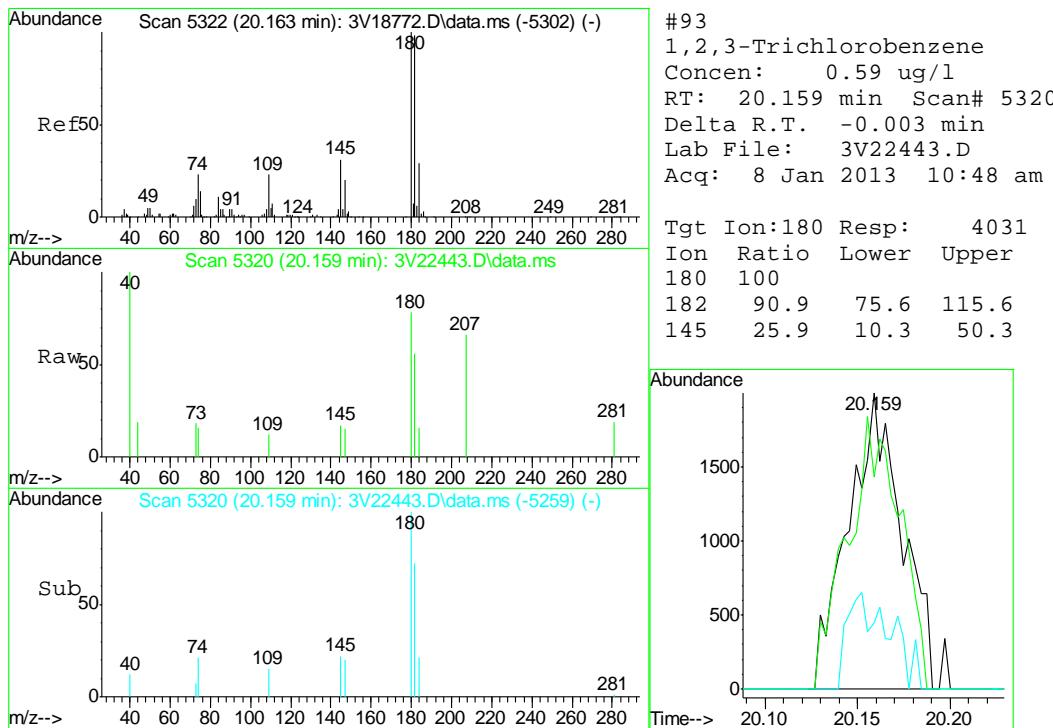
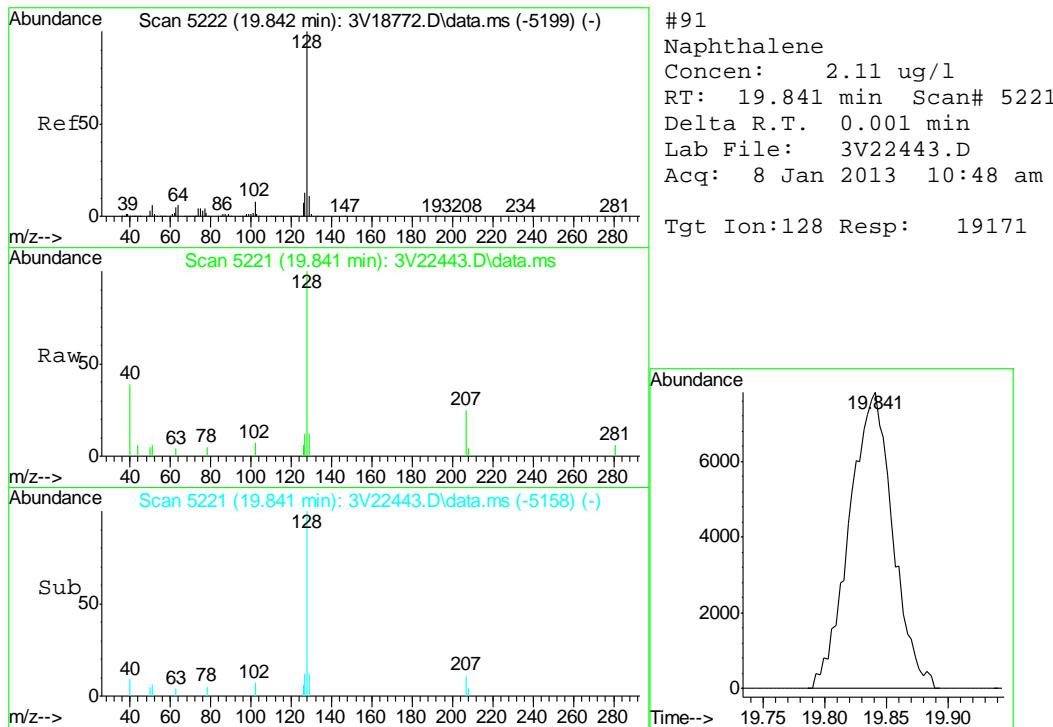


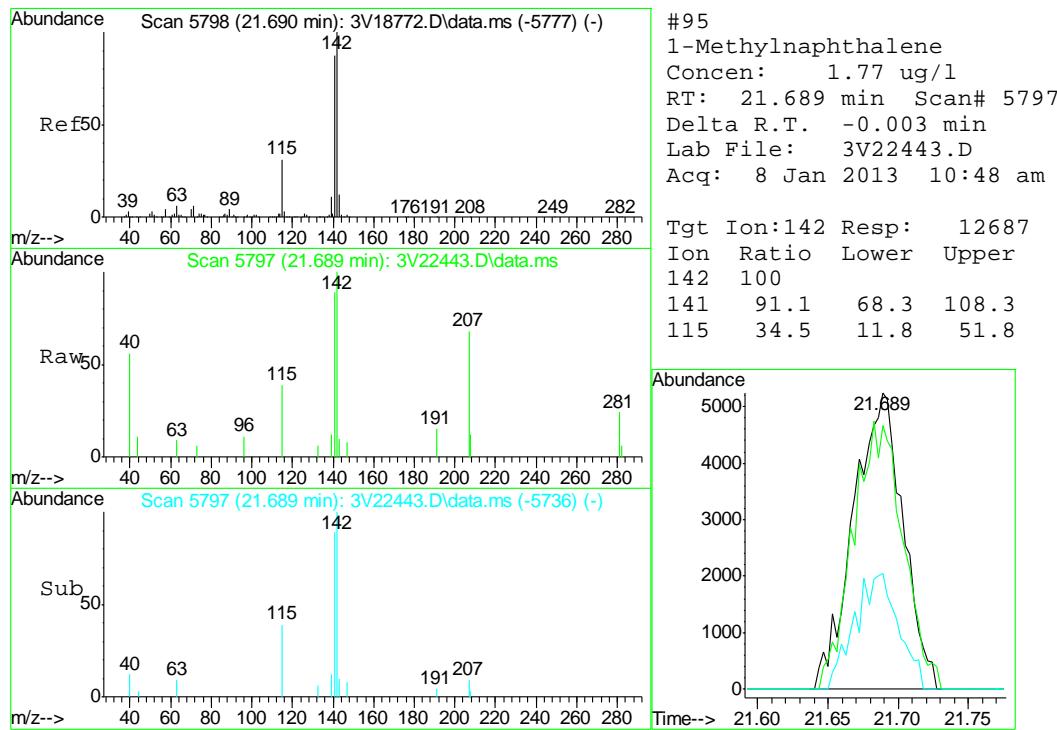
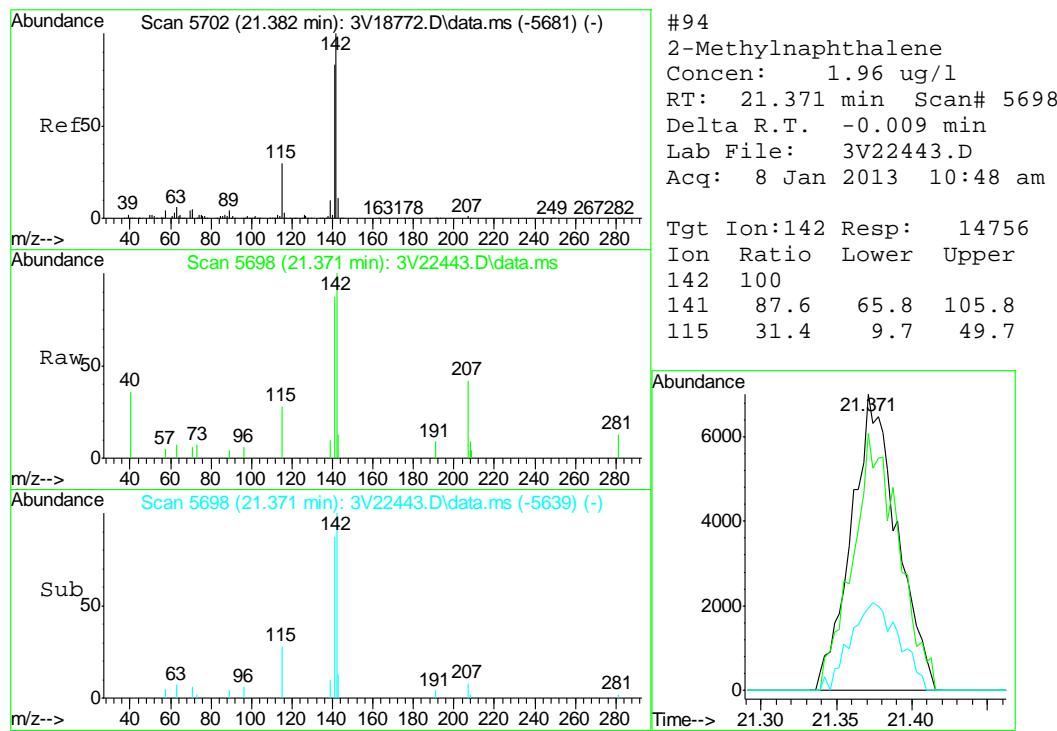














## GC/MS Semi-volatiles

---

### QC Data Summaries

∞

---

Includes the following where applicable:

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

**Method Blank Summary**

**Job Number:** D42316  
**Account:** XTOKWR XTO Energy  
**Project:** XTO Love Ranch 8

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
OP7200-MB	3G12911.D	1	01/10/13	DC	01/09/13	OP7200	E3G618

The QC reported here applies to the following samples:

**Method:** SW846 8270C BY SIM

D42316-1

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

CAS No.	Surrogate Recoveries	Limits
4165-60-0	Nitrobenzene-d5	105%
321-60-8	2-Fluorobiphenyl	101%
1718-51-0	Terphenyl-d14	160% * a
		18-150%

(a) Outside of control limits. Since the bias is high and the method blank is ND for target analytes, no further action is required.

## Blank Spike Summary

Page 1 of 1

Job Number: D42316

Account: XTOKWR XTO Energy

Project: XTO Love Ranch 8

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
OP7200-BS	3G12910.D	1	01/10/13	DC	01/09/13	OP7200	E3G618

The QC reported here applies to the following samples:

Method: SW846 8270C BY SIM

D42316-1

CAS No.	Compound	Spike ug/kg	BSP ug/kg	BSP %	Limits
83-32-9	Acenaphthene	83.3	84.9	102	68-130
120-12-7	Anthracene	83.3	73.4	88	67-130
56-55-3	Benzo(a)anthracene	83.3	61.0	73	65-130
205-99-2	Benzo(b)fluoranthene	83.3	73.0	88	44-130
207-08-9	Benzo(k)fluoranthene	83.3	72.0	86	56-131
50-32-8	Benzo(a)pyrene	83.3	76.2	91	62-130
218-01-9	Chrysene	83.3	79.5	95	70-130
53-70-3	Dibenzo(a,h)anthracene	83.3	53.8	65	55-130
206-44-0	Fluoranthene	83.3	64.8	78	70-130
86-73-7	Fluorene	83.3	70.2	84	70-130
193-39-5	Indeno(1,2,3-cd)pyrene	83.3	57.5	69	56-130
91-20-3	Naphthalene	83.3	74.2	89	70-130
129-00-0	Pyrene	83.3	79.4	95	70-130

CAS No.	Surrogate Recoveries	BSP	Limits
4165-60-0	Nitrobenzene-d5	101%	10-159%
321-60-8	2-Fluorobiphenyl	98%	19-131%
1718-51-0	Terphenyl-d14	130%	18-150%

\* = Outside of Control Limits.

# Matrix Spike/Matrix Spike Duplicate Summary

Page 1 of 1

Job Number: D42316

Account: XTOKWR XTO Energy

Project: XTO Love Ranch 8

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
OP7200-MS	3G12915.D	4	01/10/13	DC	01/09/13	OP7200	E3G618
OP7200-MSD	3G12914.D	4	01/10/13	DC	01/09/13	OP7200	E3G618
D42316-1	3G12922.D	1	01/10/13	DC	01/09/13	OP7200	E3G618
D42316-1	3G12912.D	4	01/10/13	DC	01/09/13	OP7200	E3G618

The QC reported here applies to the following samples:

Method: SW846 8270C BY SIM

D42316-1

CAS No.	Compound	D42316-1 ug/kg	Spike ug/kg	MS ug/kg	MS %	MSD ug/kg	MSD %	RPD	Limits Rec/RPD
83-32-9	Acenaphthene	ND	131	141	107	215	163* <sup>a</sup>	42* <sup>b</sup>	25-151/30
120-12-7	Anthracene	ND	131	164	125	168	128	2	39-159/30
56-55-3	Benzo(a)anthracene	ND	131	122	93	127	97	4	39-168/30
205-99-2	Benzo(b)fluoranthene	ND	131	133	101	129	98	3	24-163/30
207-08-9	Benzo(k)fluoranthene	ND	131	94.4	72	111	84	16	10-188/30
50-32-8	Benzo(a)pyrene	ND	131	107	81	113	86	5	32-144/30
218-01-9	Chrysene	ND	131	161	123	168	128	4	43-150/30
53-70-3	Dibenzo(a,h)anthracene	ND	131	89.6	68	90.5	69	1	21-152/30
206-44-0	Fluoranthene	ND	131	173	132	178	135	3	36-157/30
86-73-7	Fluorene	437	131	845	311* <sup>c</sup>	829	298* <sup>c</sup>	2	10-182/30
193-39-5	Indeno(1,2,3-cd)pyrene	ND	131	87.8	67	89.5	68	2	20-154/30
91-20-3	Naphthalene	1190 <sup>d</sup>	131	1480	221* <sup>c</sup>	1510	243* <sup>c</sup>	2	10-163/30
129-00-0	Pyrene	80.0	131	216	104	227	112	5	25-180/30

CAS No.	Surrogate Recoveries	MS	MSD	D42316-1	D42316-1	Limits
4165-60-0	Nitrobenzene-d5	34%	42%	84%	42%	10-159%
321-60-8	2-Fluorobiphenyl	52%	70%	50%	48%	19-131%
1718-51-0	Terphenyl-d14	104%	110%	73%	107%	18-150%

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

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

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

(d) Result is from Run #2.

\* = Outside of Control Limits.

8.3.1  
8



## GC/MS Semi-volatiles

---

Raw Data

---

## Quantitation Report (QT Reviewed)

Manual Integrations  
APPROVED  
(compounds with "m" flag)

Judy Nelson  
01/11/13 11:06

Data Path : C:\msdchem\1\DATA\011013\  
Data File : 3g12912.D  
Acq On : 10 Jan 13 3:08 pm  
Operator : DONC  
Sample : D42316-1, 4x  
Misc : OP7200,E3G618,30.03,,,1,4  
ALS Vial : 6 Sample Multiplier: 1

Quant Time: Jan 11 09:02:28 2013  
Quant Method : C:\msdchem\1\METHODS\SIMPE3G611.M  
Quant Title : PAHSIM BASE  
QLast Update : Thu Jan 10 14:18:35 2013  
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Naphthalene-d8	5.621	136	135536	4.0000	ug/mL	0.00
6) Acenaphthene-d10	7.337	164	92581	4.0000	ug/mL	0.01
15) Phenanthrene-d10	8.819	188	129785	4.0000	ug/mL	0.00
19) Chrysene-d12	11.450	240	89785	4.0000	ug/mL	0.00
24) Perylene-d12	12.810	264	68780	4.0000	ug/mL	0.00

## System Monitoring Compounds

2) Nitrobenzene-d5	4.935	82	63983m	5.2483	ug/mL	-0.01
Spiked Amount	50.000	Range	25 - 135	Recovery	=	10.50%#
7) 2-Fluorobiphenyl	6.676	172	236998	6.0253	ug/mL	0.00
Spiked Amount	50.000	Range	25 - 135	Recovery	=	12.06%#
21) Terphenyl-d14	10.402	244	163950	13.4198	ug/mL	0.00
Spiked Amount	50.000	Range	25 - 135	Recovery	=	26.84%

## Target Compounds

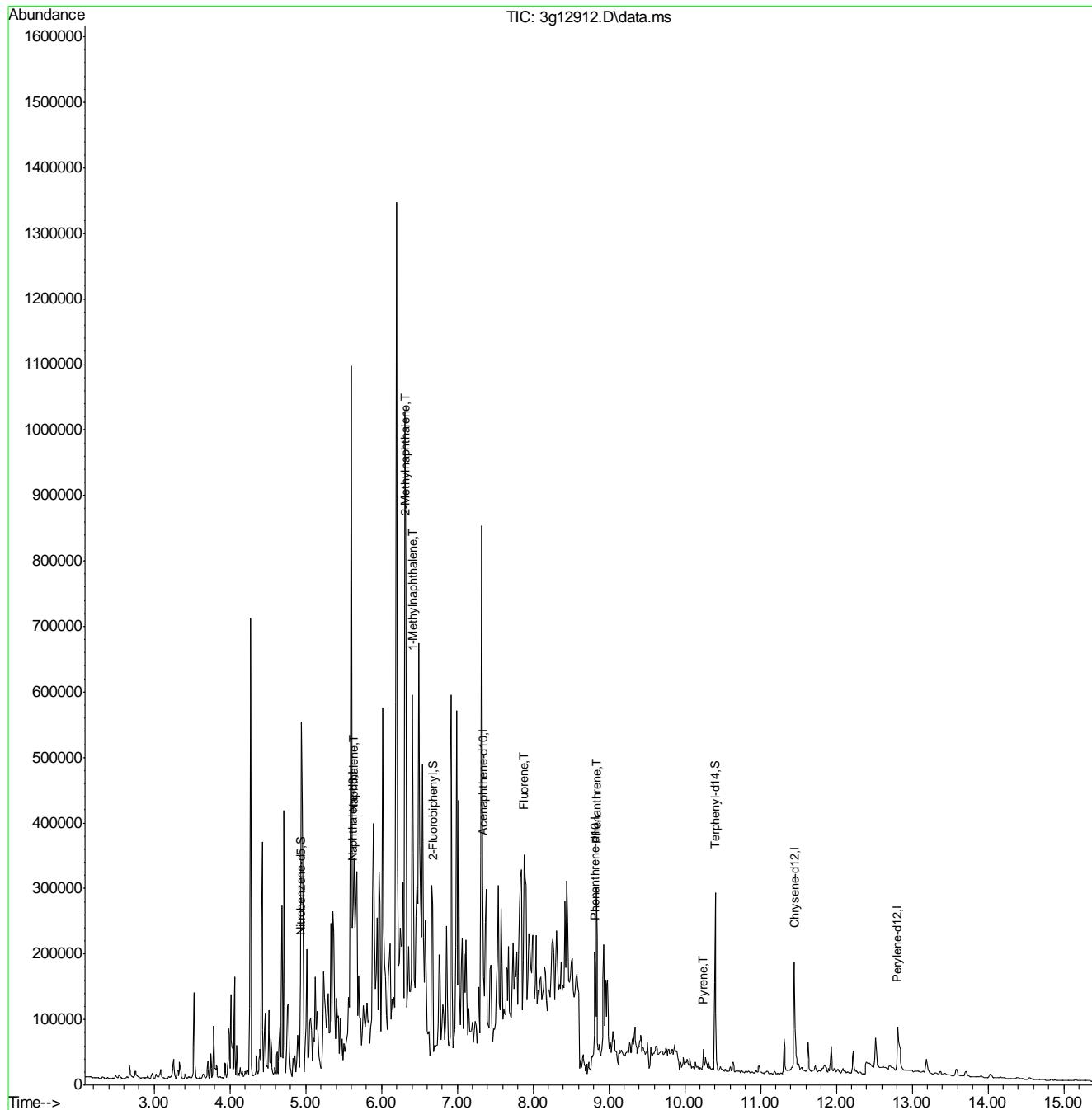
				Qvalue
3) N-Nitrosodimethylamine	2.363	74	35	N.D.
4) N-Nitrosodi-propylamine	0.000	70	0	N.D. d
5) Naphthalene	5.633	128	220836	5.6665 ug/mL 90
8) 2-Methylnaphthalene	6.319	142	538416	18.0980 ug/mL 95
9) 1-Methylnaphthalene	6.406	142	184588m	7.0939 ug/mL
10) Acenaphthylene	0.000	152	0	N.D. d
11) Acenaphthene	0.000	154	0	N.D. d
12) Dibenzofuran	0.000	168	0	N.D. d
13) Fluorene	7.881	166	97530m	2.6190 ug/mL
14) Diphenylamine	0.000	169	0	N.D. d
16) Phenanthrene	8.835	178	186818	3.7249 ug/mL 68
17) Anthracene	0.000	178	0	N.D. d
18) Fluoranthene	0.000	202	0	N.D. d
20) Pyrene	10.244	202	20757	0.4329 ug/mL# 40
22) Benzo(a)anthracene	0.000	228	0	N.D. d
23) Chrysene	0.000	228	0	N.D. d
25) Benzo(b)fluoranthene	0.000	252	0	N.D. d
26) Benzo(k)fluoranthene	0.000	252	0	N.D. d
27) Benzo(a)pyrene	0.000	252	0	N.D. d
28) Indeno(1,2,3-cd)pyrene	0.000	276	0	N.D. d
29) Dibenz(a,h)anthracene	0.000	278	0	N.D. d
30) Benzo(g,h,i)perylene	14.345	276	805	N.D.

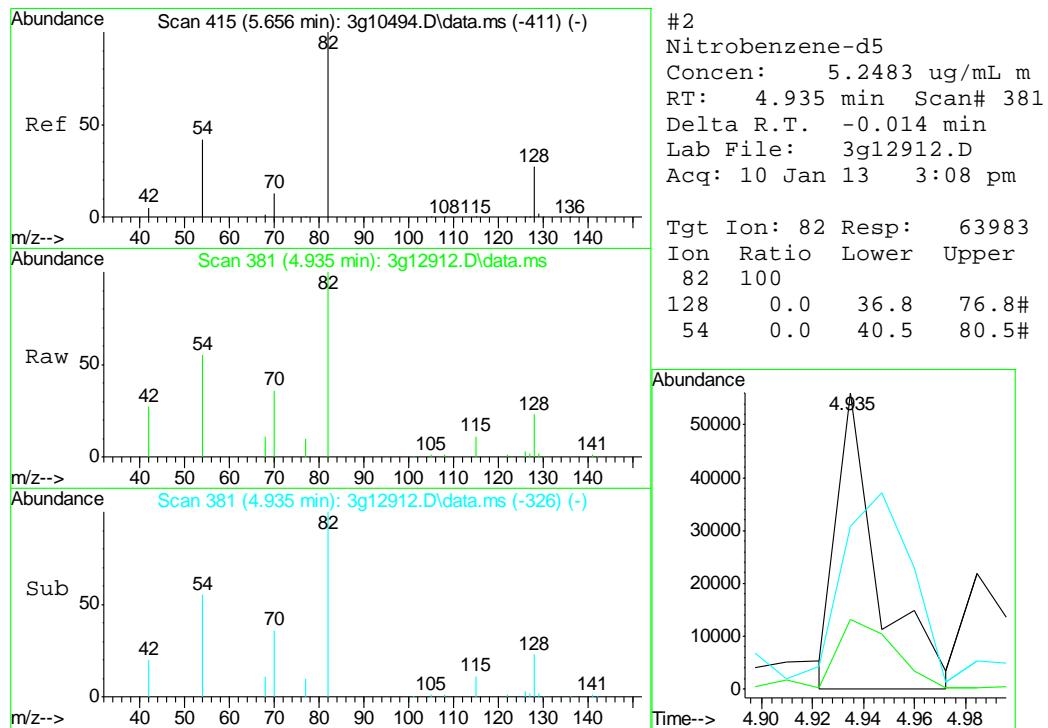
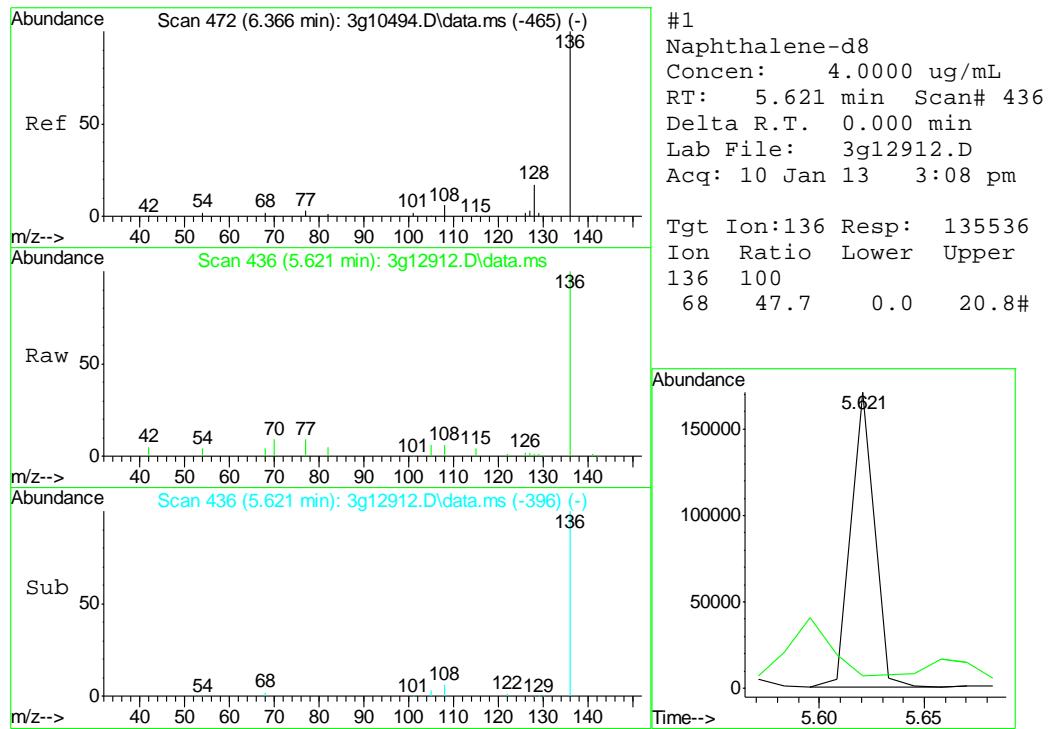
(#) = qualifier out of range (m) = manual integration (+) = signals summed

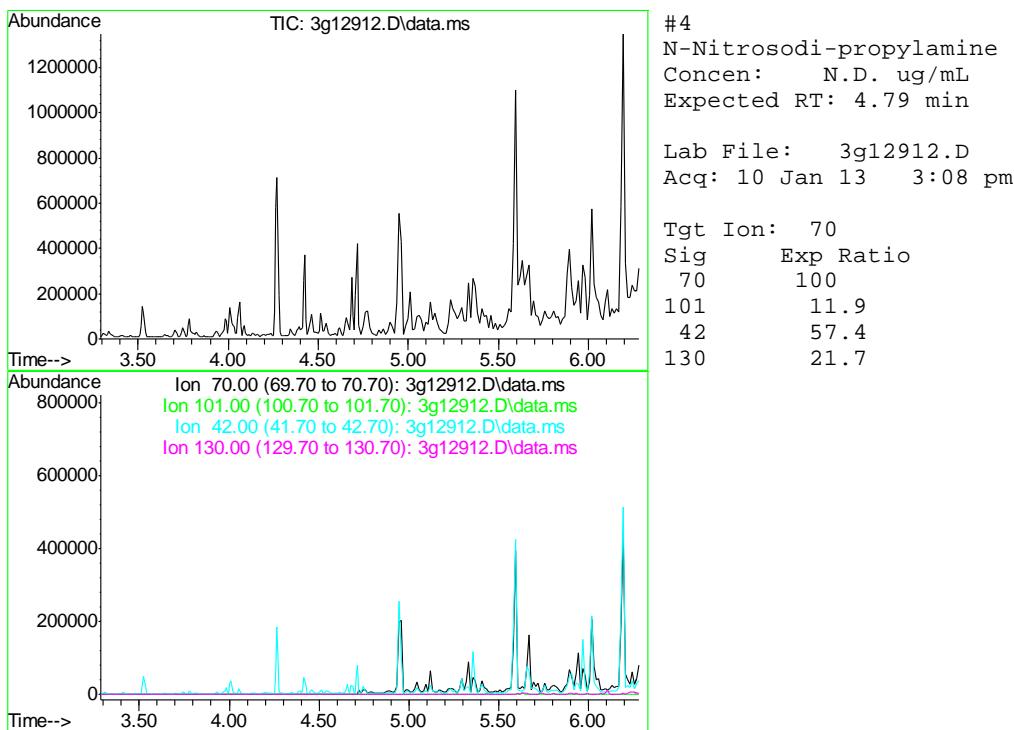
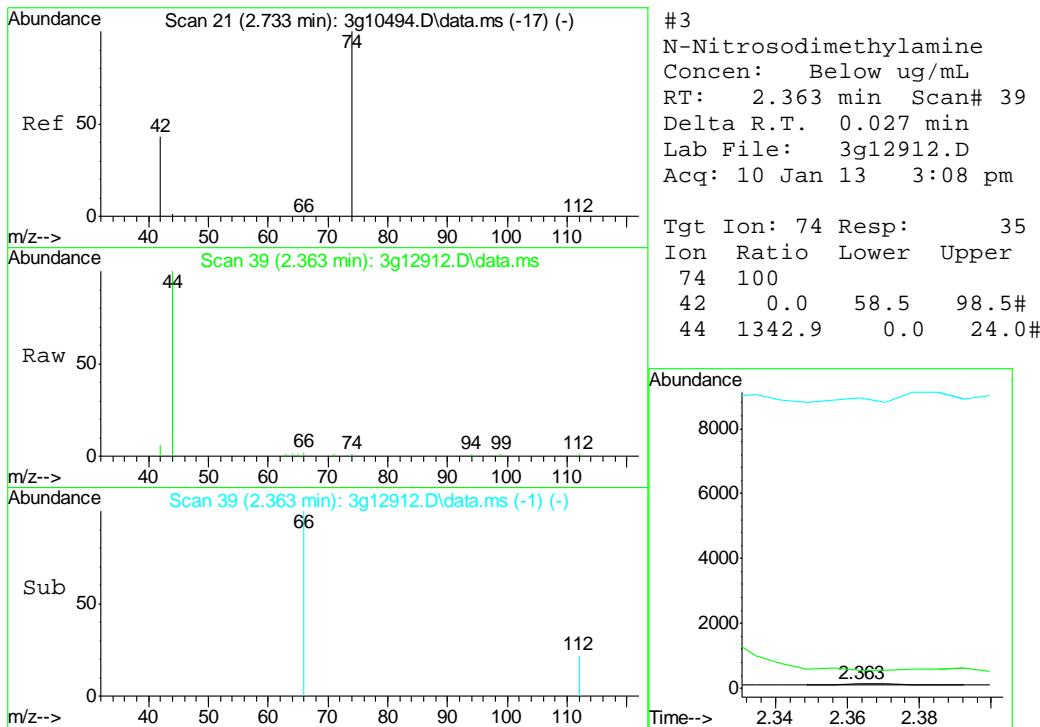
## Quantitation Report (QT Reviewed)

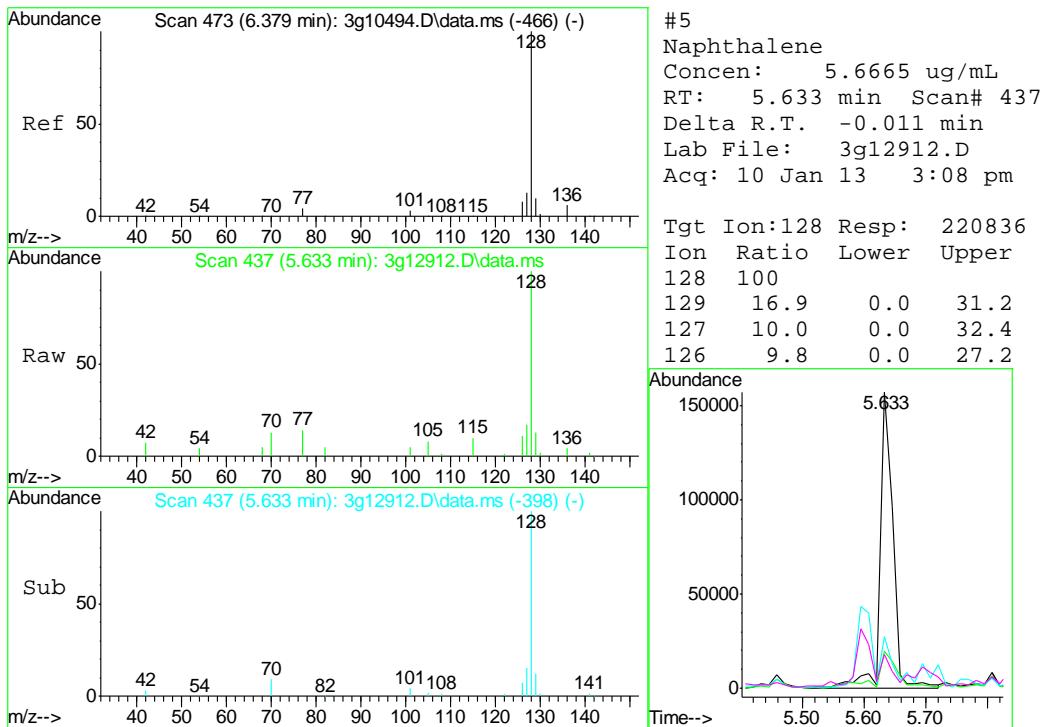
Data Path : C:\msdchem\1\DATA\011013\  
 Data File : 3g12912.D  
 Acq On : 10 Jan 13 3:08 pm  
 Operator : DONC  
 Sample : D42316-1, 4x  
 Misc : OP7200,E3G618,30.03,,,1,4  
 ALS Vial : 6 Sample Multiplier: 1

Quant Time: Jan 11 09:02:28 2013  
 Quant Method : C:\msdchem\1\METHODS\SIMPE3G611.M  
 Quant Title : PAHSIM BASE  
 QLast Update : Thu Jan 10 14:18:35 2013  
 Response via : Initial Calibration



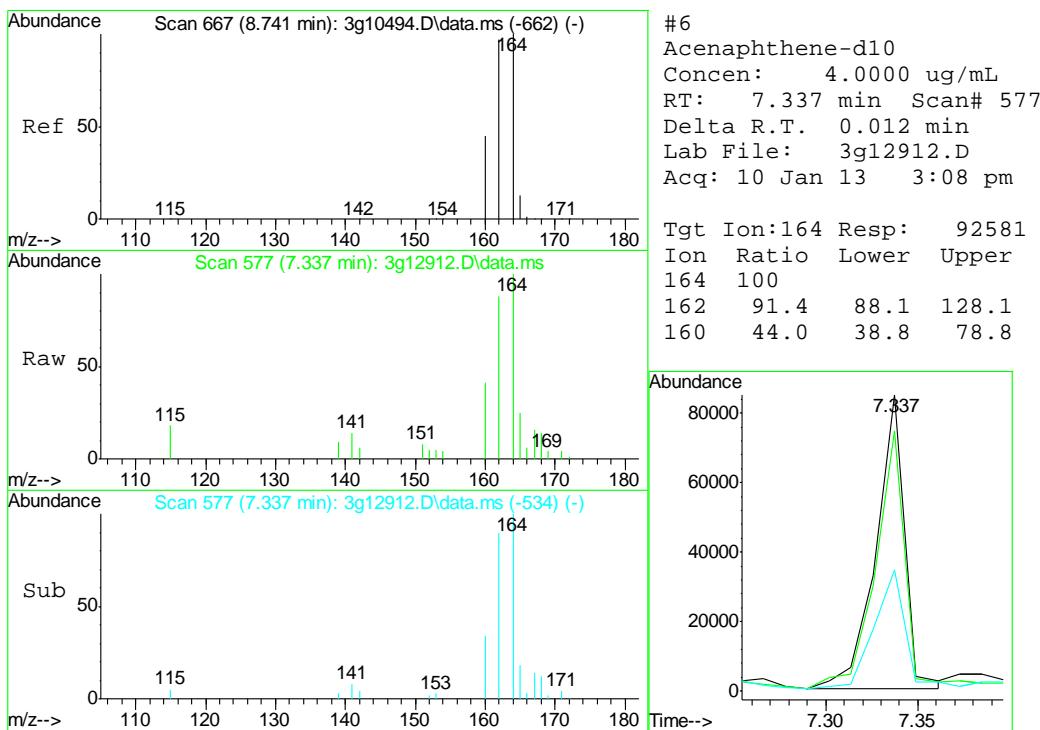


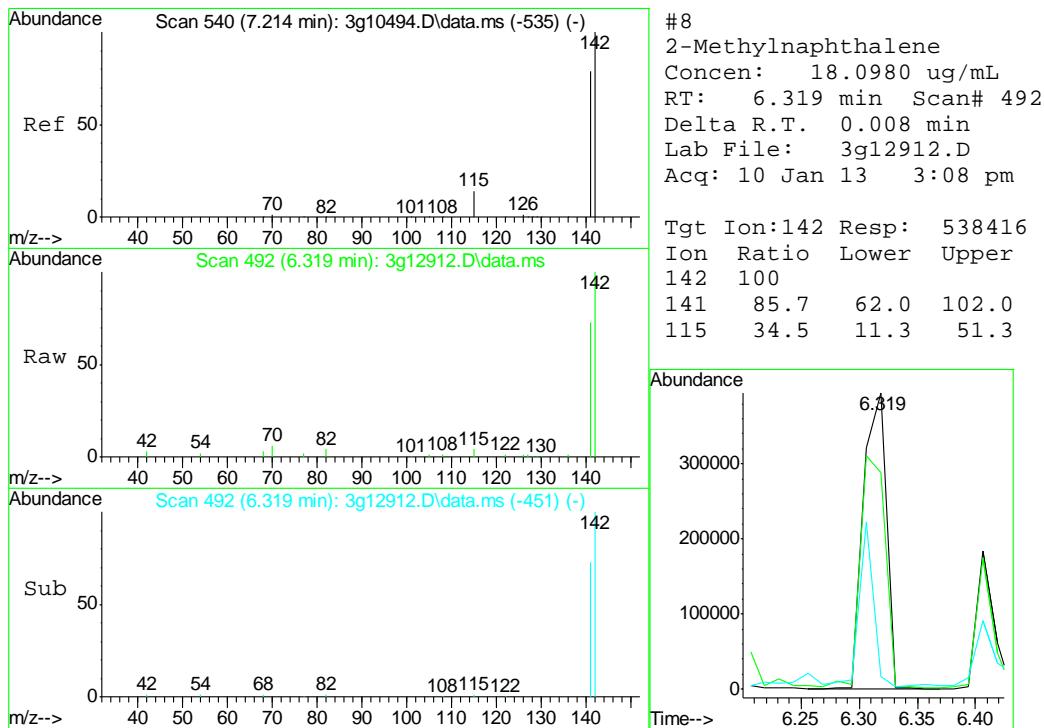
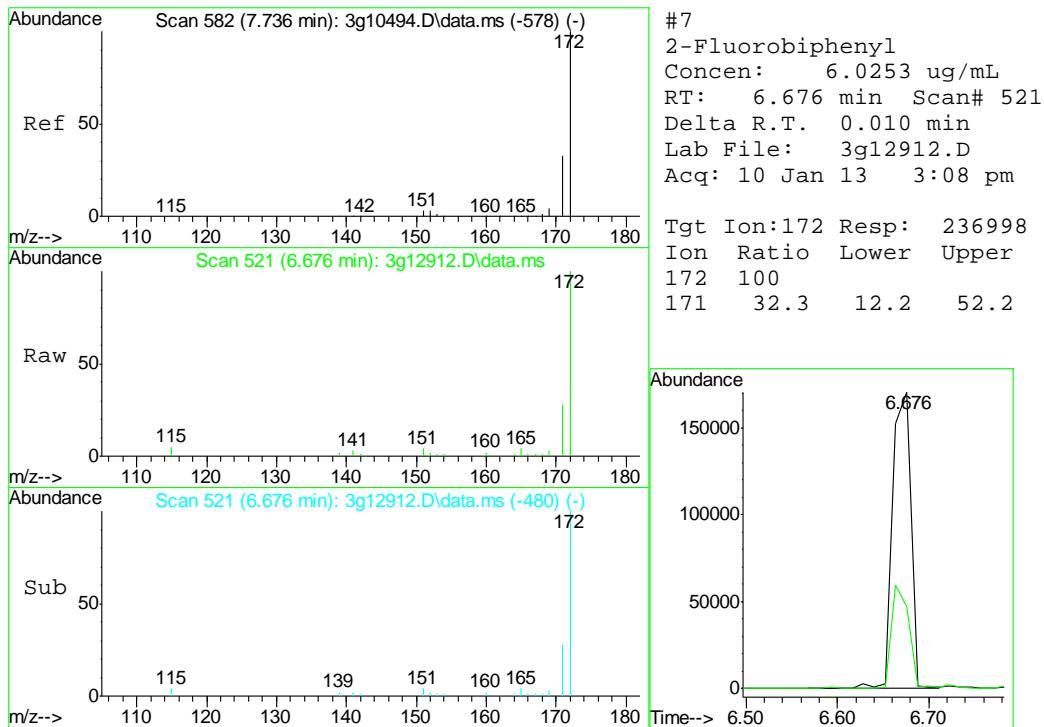


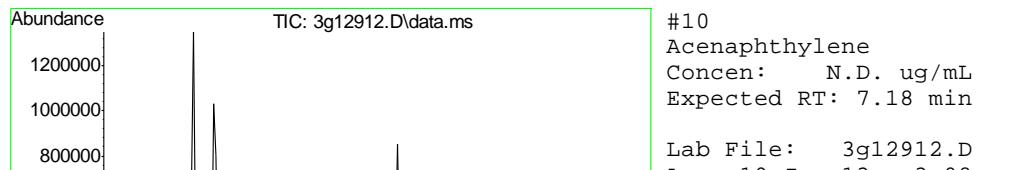
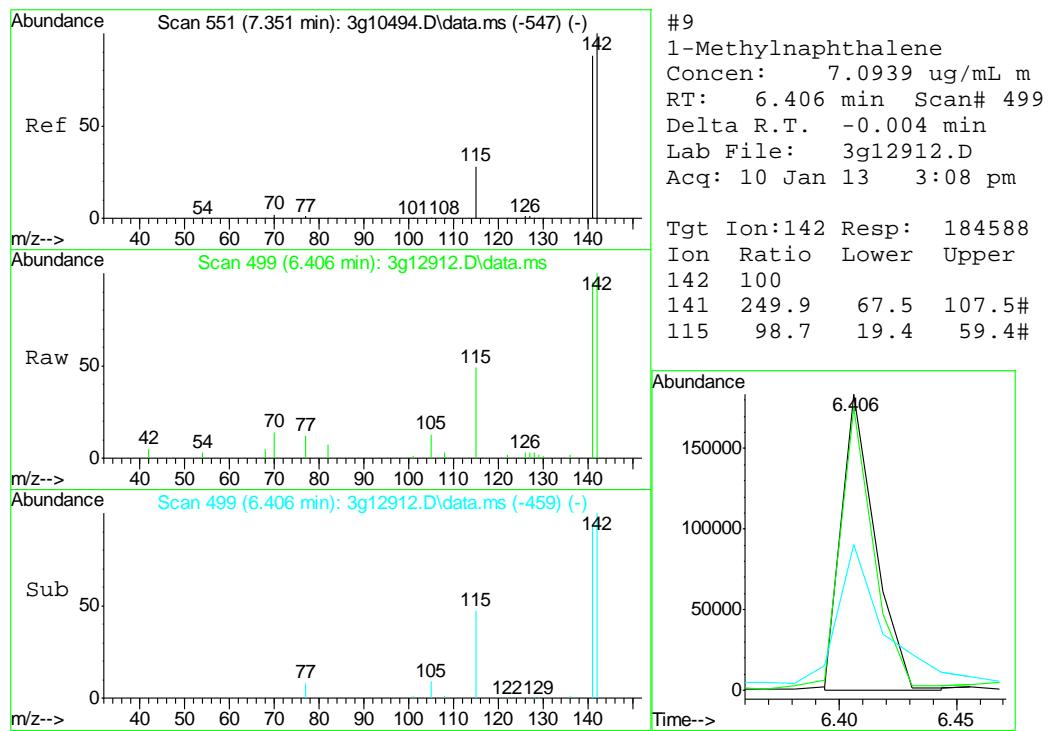


9.1.1

6

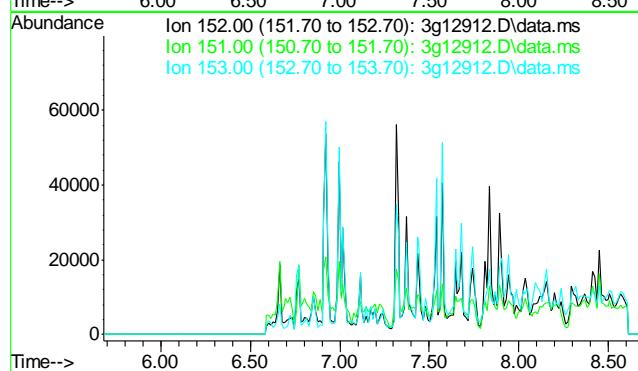


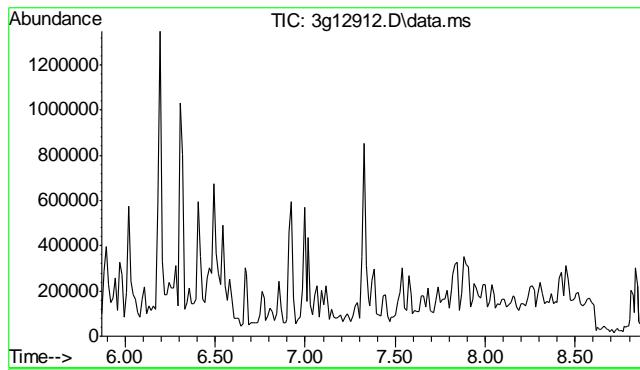




Lab File: 3g12912.D  
Acq: 10 Jan 13 3:08 pm

Tgt Ion: 152  
Sig Exp Ratio  
152 100  
151 19.2  
153 12.9

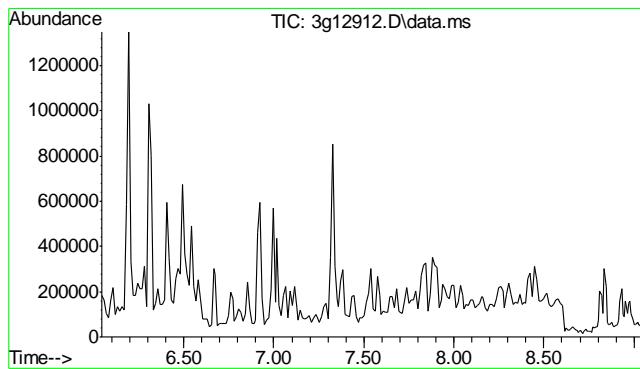
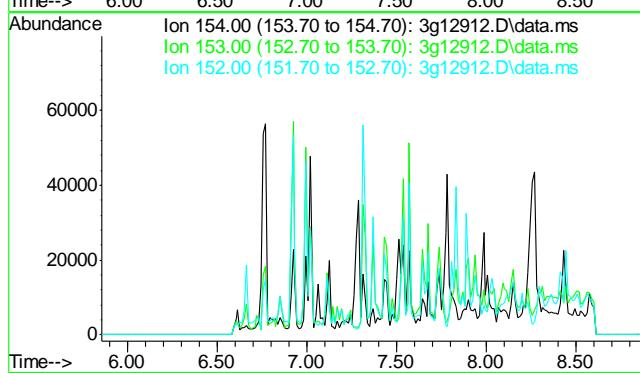




#11  
 Acenaphthene  
 Concen: N.D. ug/mL  
 Expected RT: 7.36 min

Lab File: 3g12912.D  
 Acq: 10 Jan 13 3:08 pm

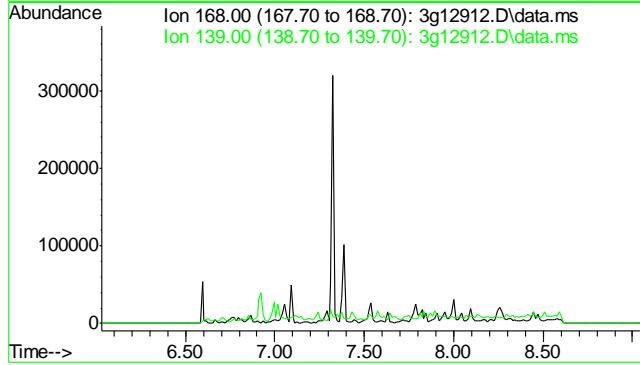
Tgt Ion:	Sig	Exp Ratio
154	100	
153	102.4	
152	50.0	

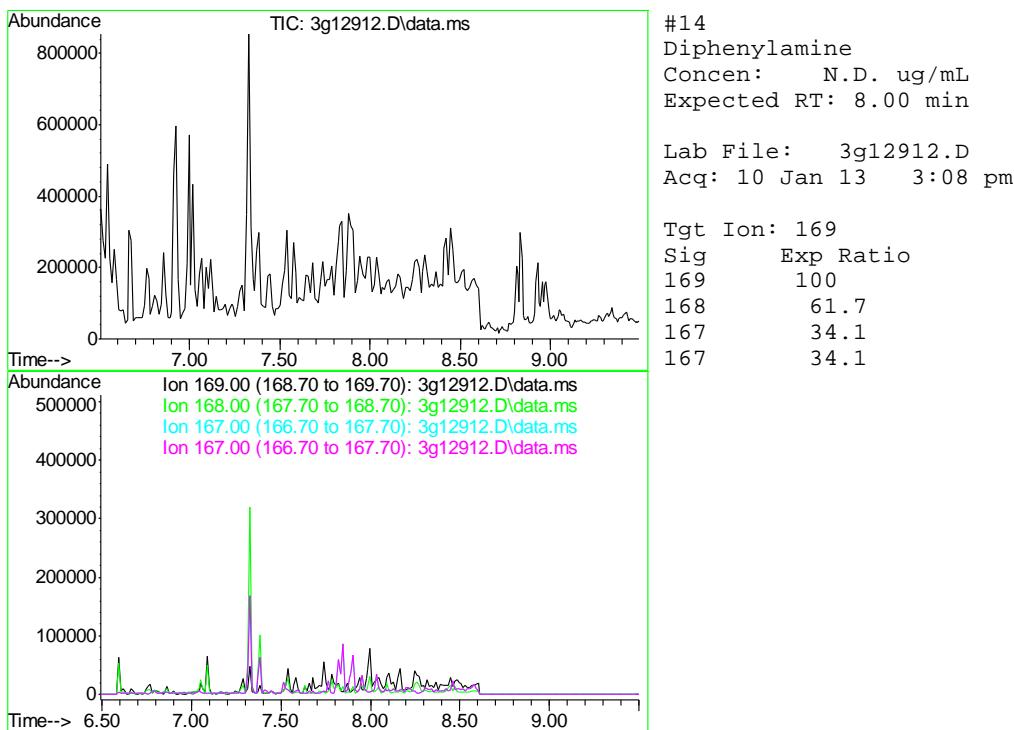
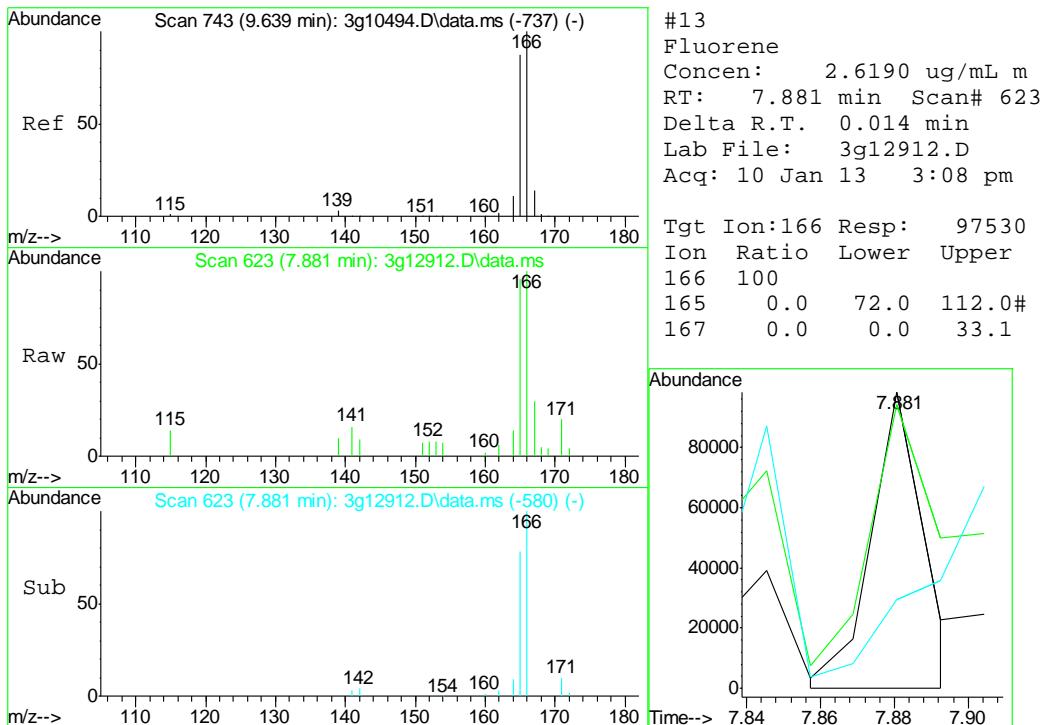


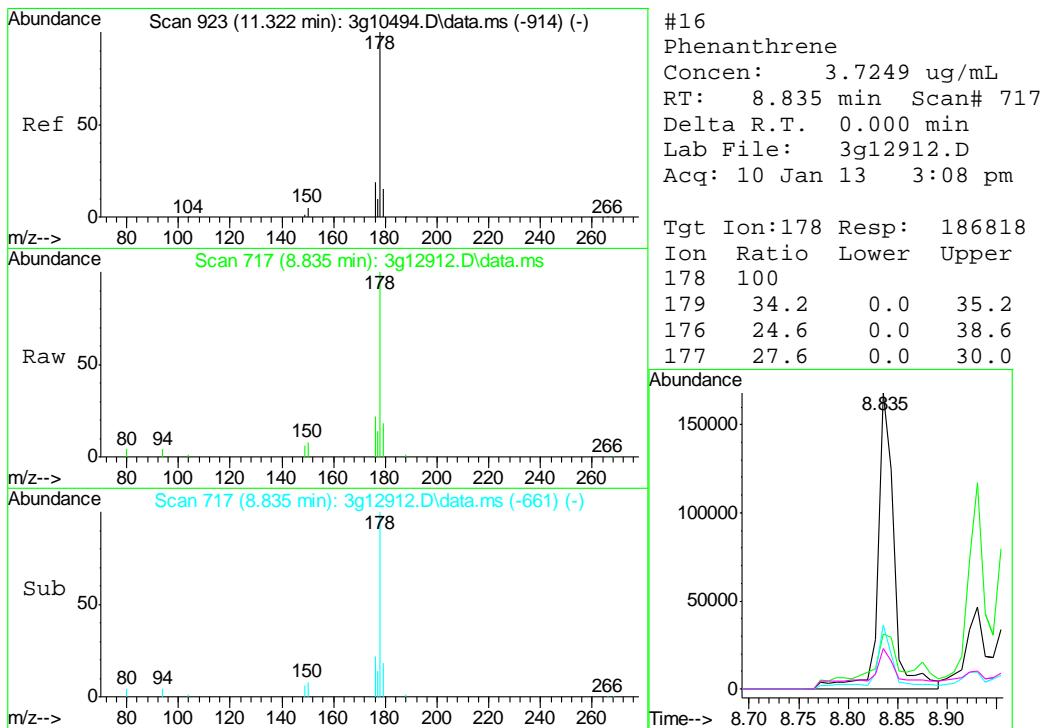
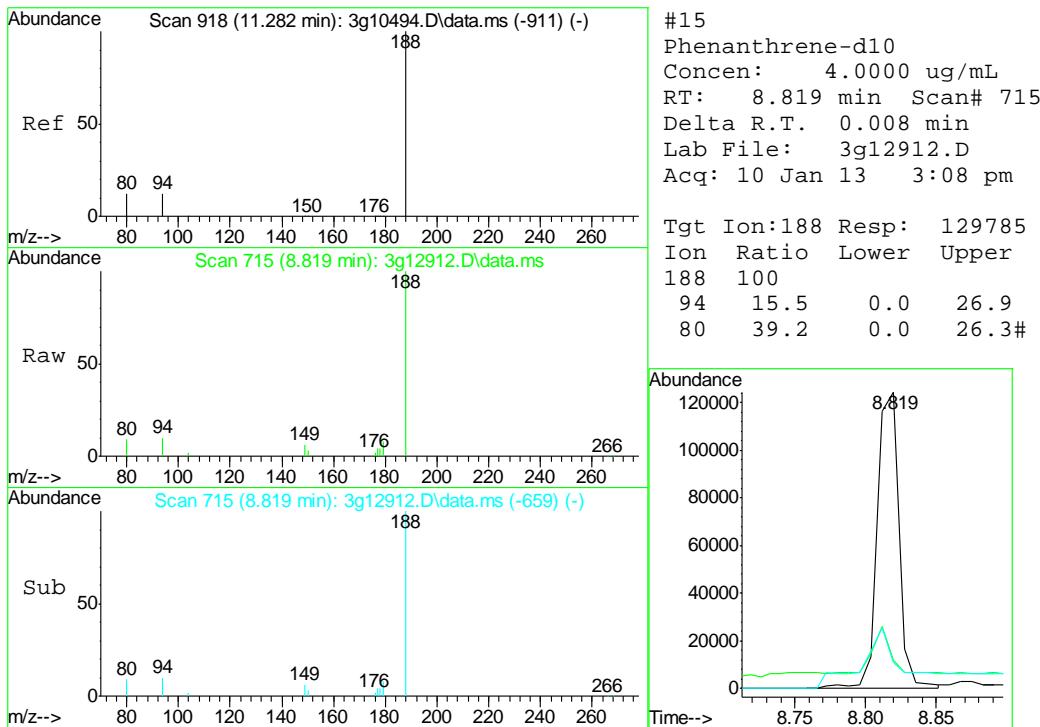
#12  
 Dibenzofuran  
 Concen: N.D. ug/mL  
 Expected RT: 7.54 min

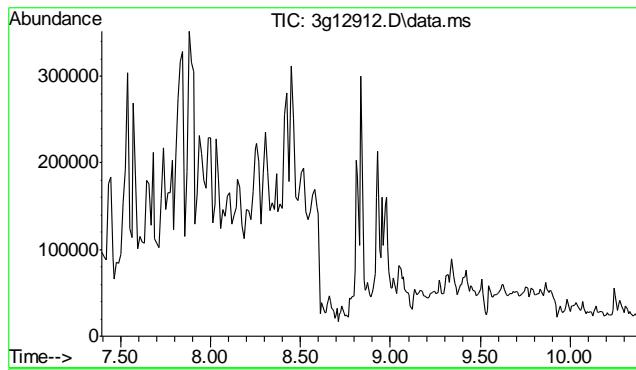
Lab File: 3g12912.D  
 Acq: 10 Jan 13 3:08 pm

Tgt Ion:	Sig	Exp Ratio
168	100	
139	33.4	

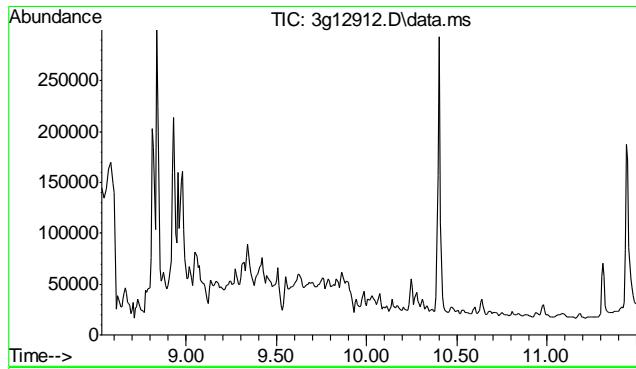
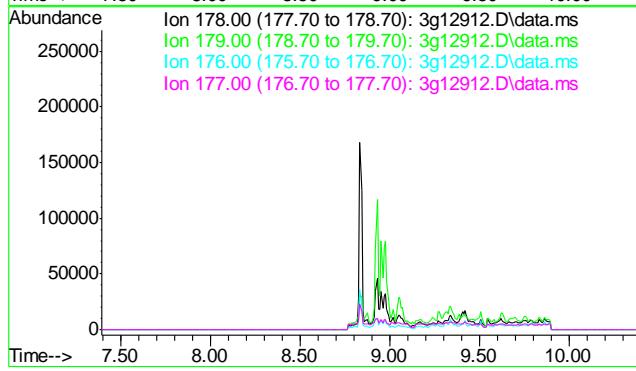




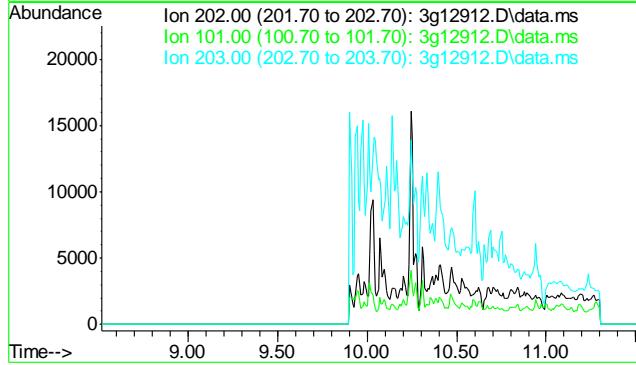


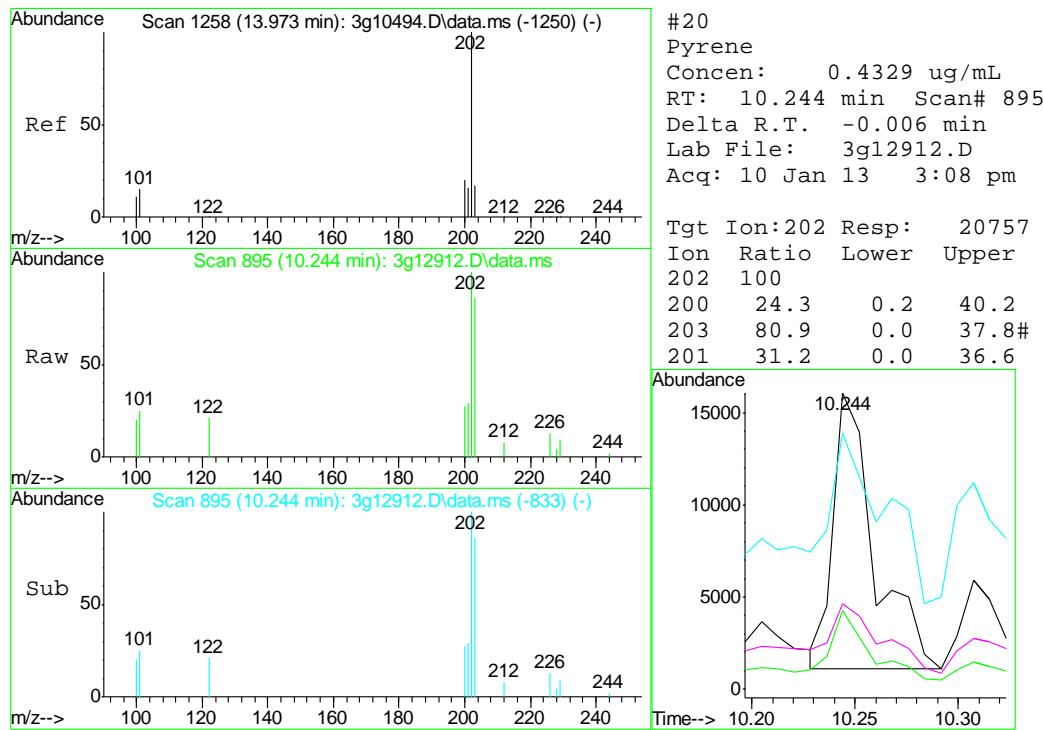
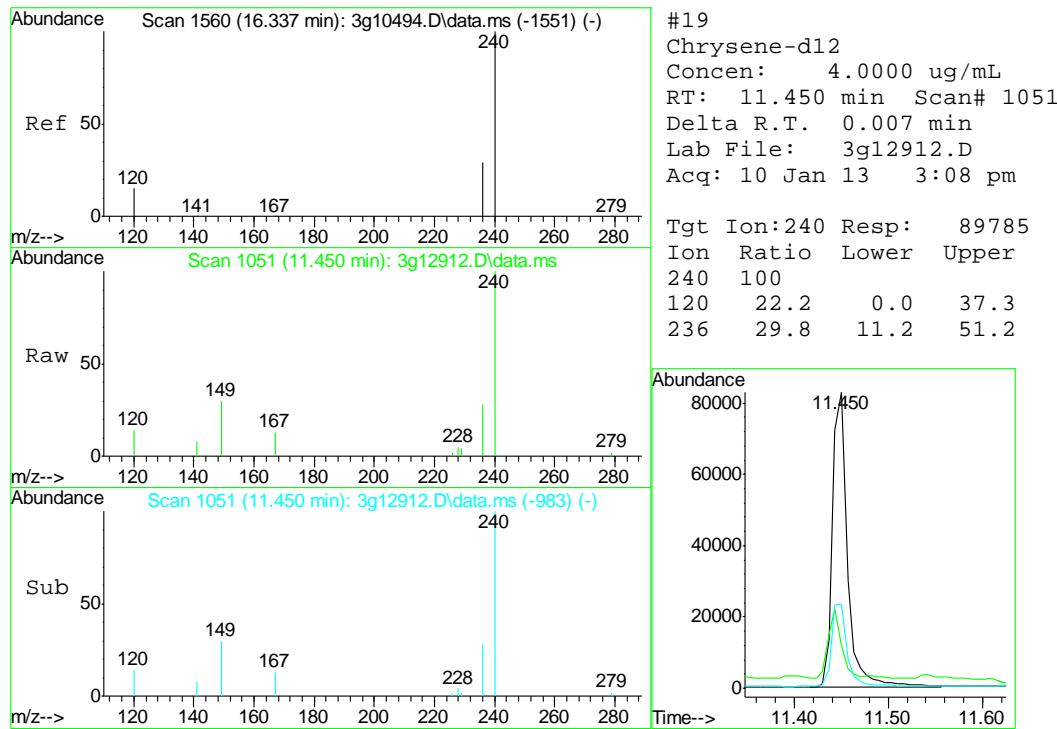


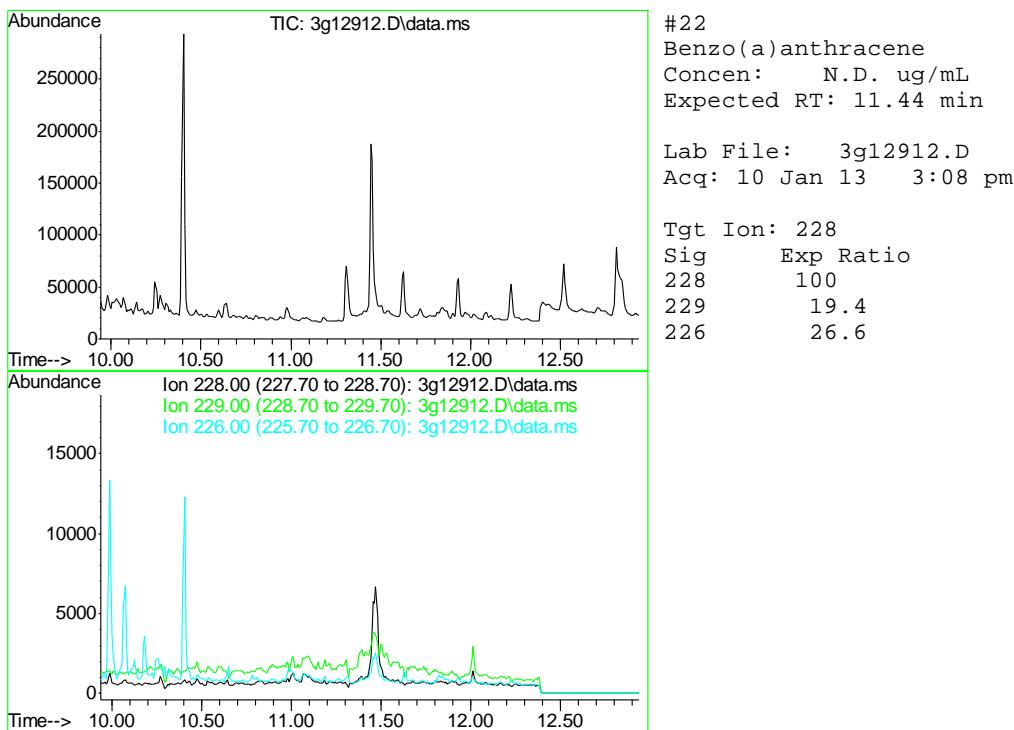
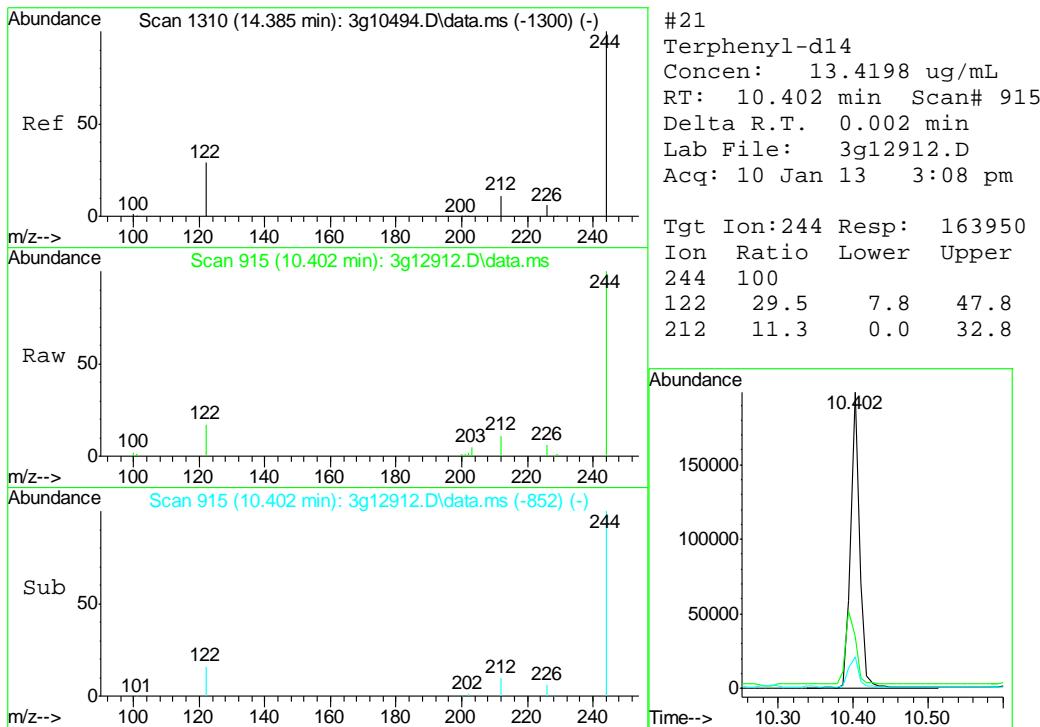
#17  
Anthracene  
Concen: N.D. ug/mL  
Expected RT: 8.89 min  
  
Lab File: 3g12912.D  
Acq: 10 Jan 13 3:08 pm  
  
Tgt Ion: 178  
Sig Exp Ratio  
178 100  
179 15.1  
176 18.2  
177 8.7

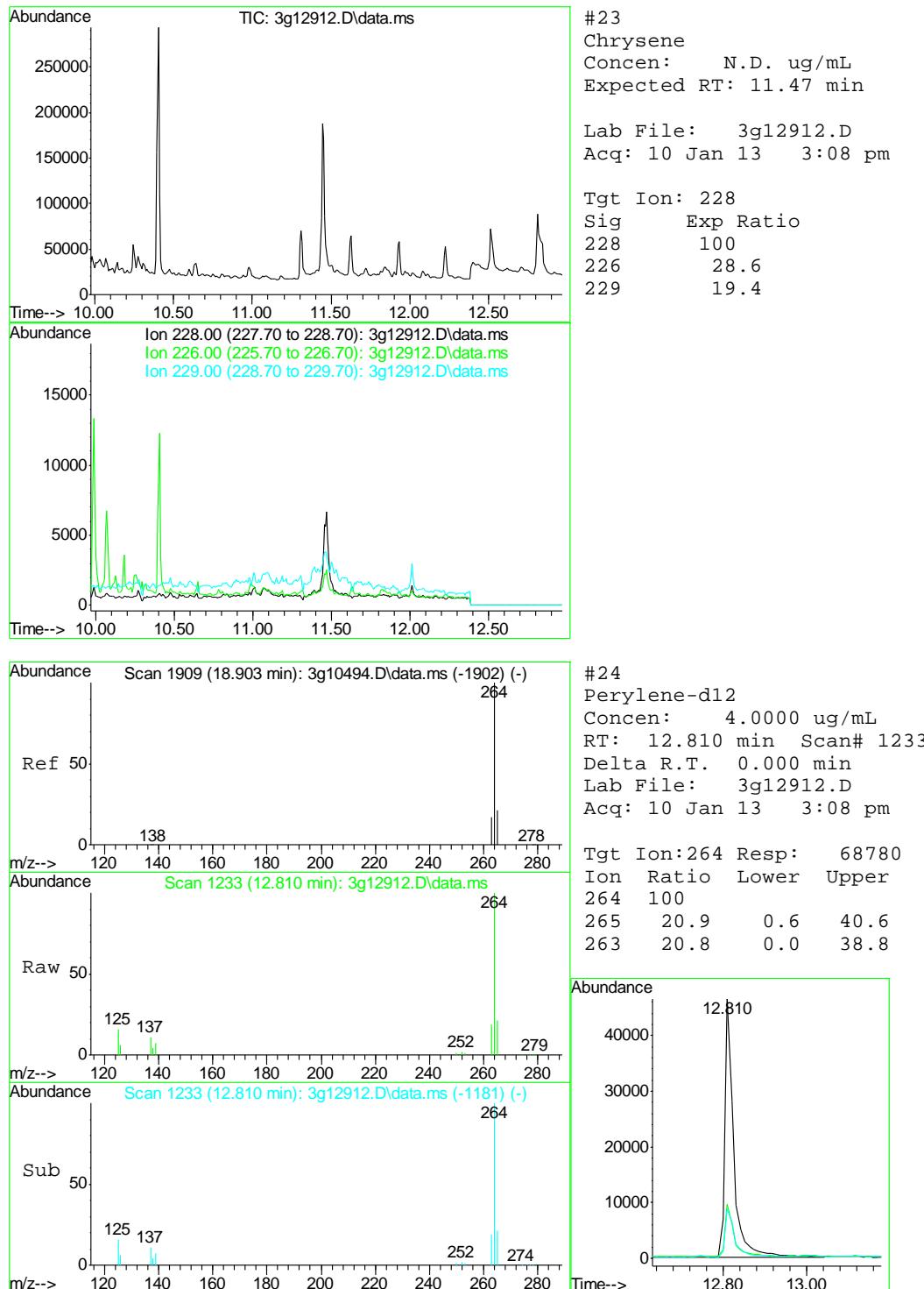


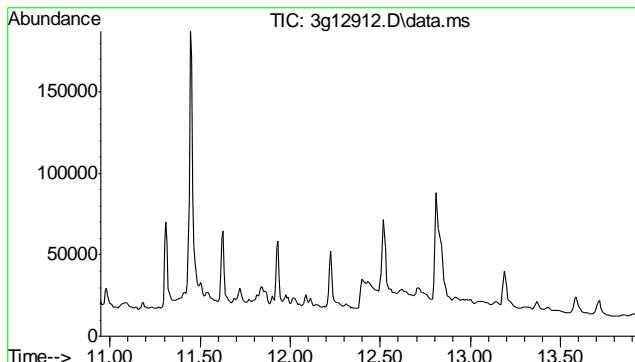
#18  
Fluoranthene  
Concen: N.D. ug/mL  
Expected RT: 10.02 min  
  
Lab File: 3g12912.D  
Acq: 10 Jan 13 3:08 pm  
  
Tgt Ion: 202  
Sig Exp Ratio  
202 100  
101 12.6  
203 17.4



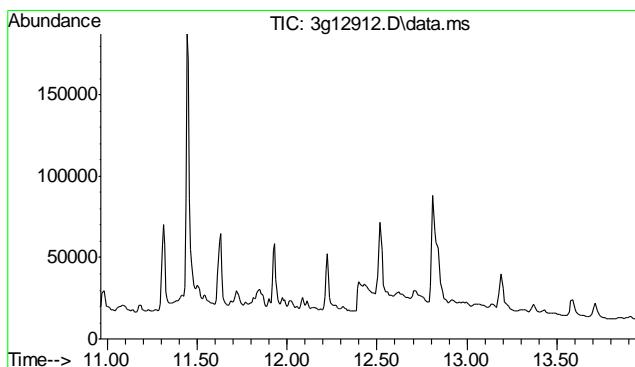
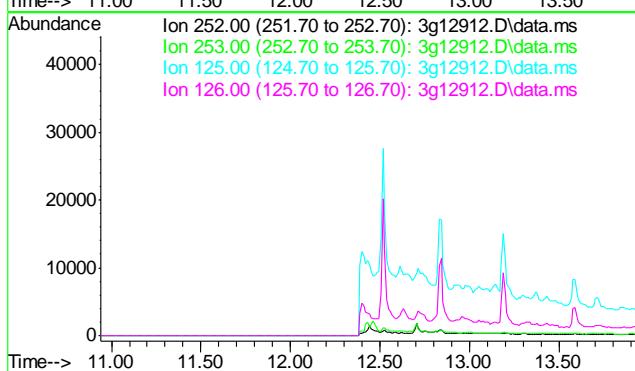




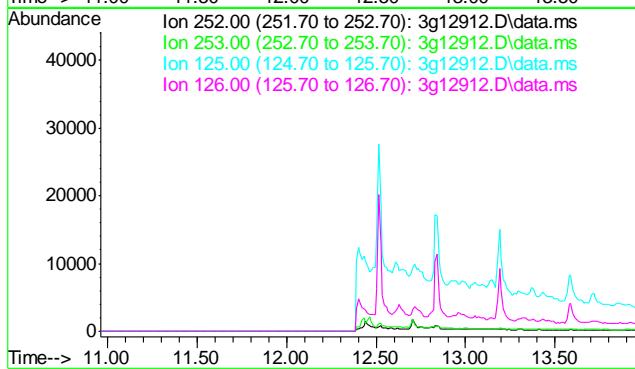


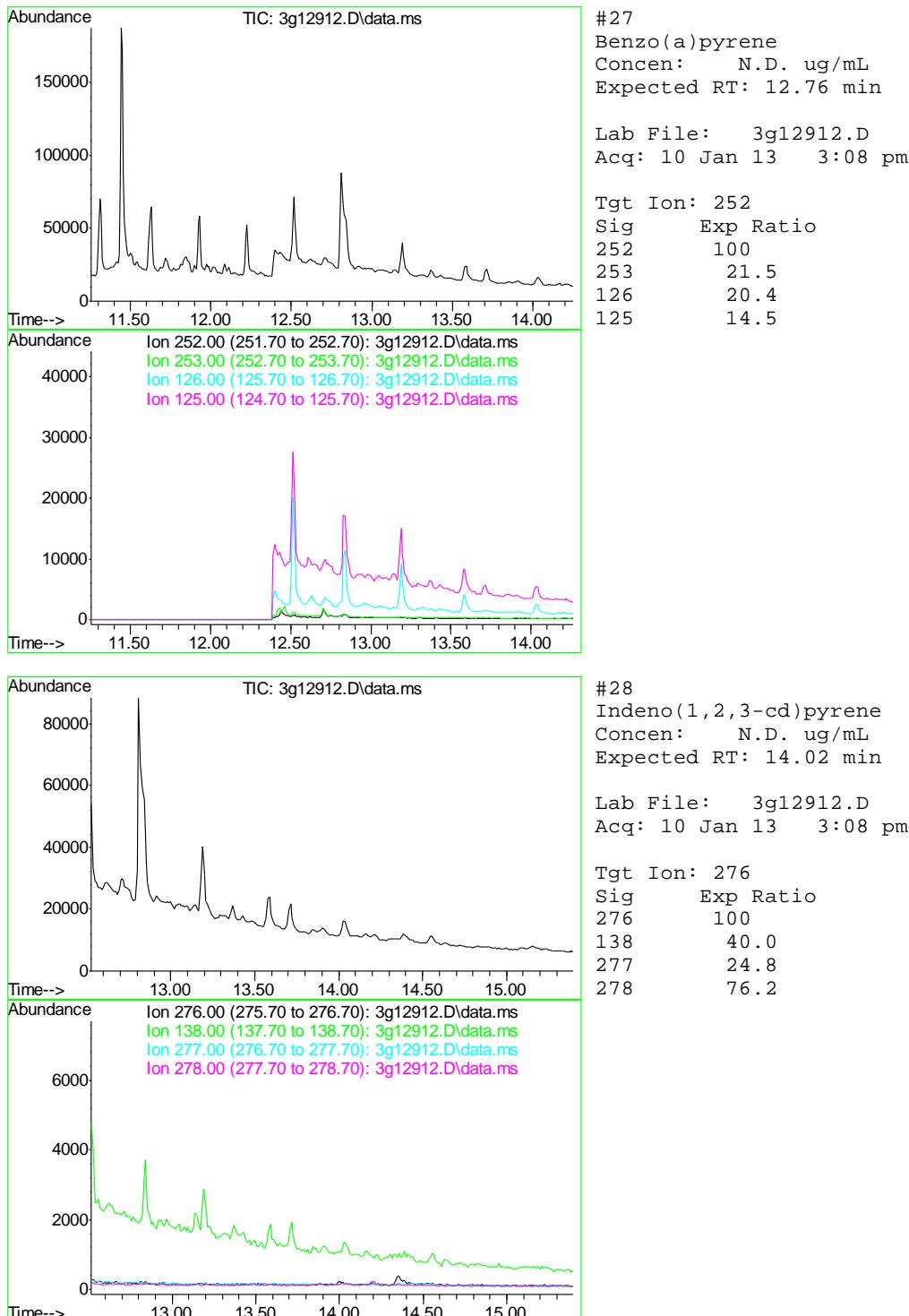


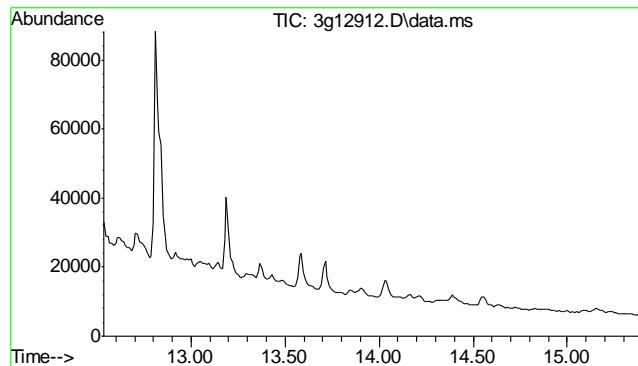
#25  
Benzo(b)fluoranthene  
Concen: N.D. ug/mL  
Expected RT: 12.44 min  
  
Lab File: 3g12912.D  
Acq: 10 Jan 13 3:08 pm  
  
Tgt Ion: 252  
Sig Exp Ratio  
252 100  
253 51.5  
125 13.2  
126 46.9



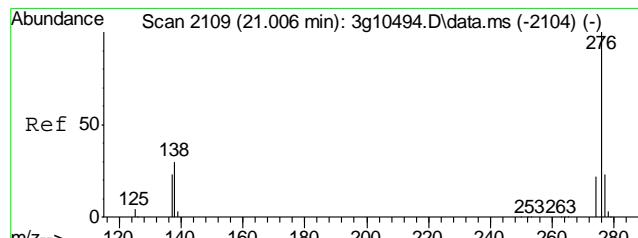
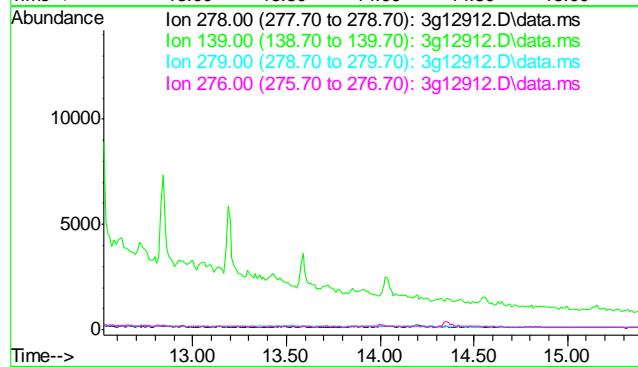
#26  
Benzo(k)fluoranthene  
Concen: N.D. ug/mL  
Expected RT: 12.46 min  
  
Lab File: 3g12912.D  
Acq: 10 Jan 13 3:08 pm  
  
Tgt Ion: 252  
Sig Exp Ratio  
252 100  
253 37.3  
125 9.6  
126 34.1





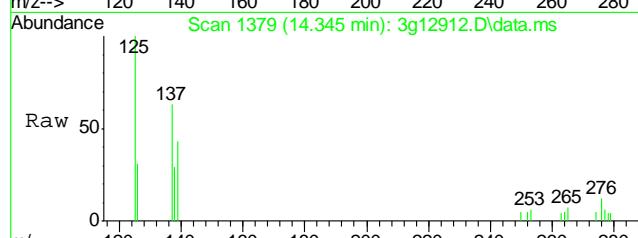


#29  
 Dibenz(a,h)anthracene  
 Concen: N.D. ug/mL  
 Expected RT: 14.03 min  
 Lab File: 3g12912.D  
 Acq: 10 Jan 13 3:08 pm  
 Tgt Ion: 278  
 Sig Exp Ratio  
 278 100  
 139 30.8  
 279 22.9  
 276 131.2

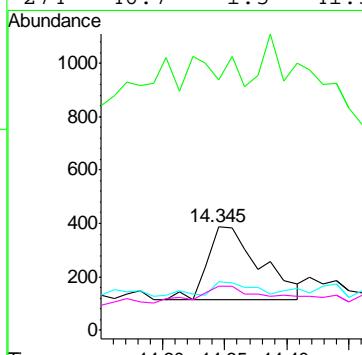
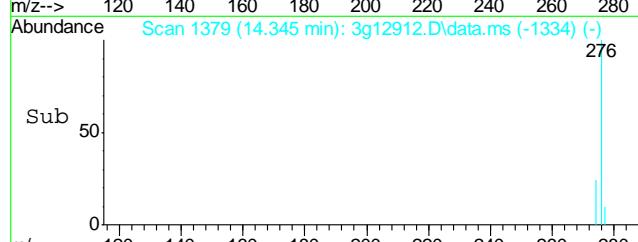


#30  
 Benzo(g,h,i)perylene  
 Concen: Below ug/mL  
 RT: 14.345 min Scan# 1379  
 Delta R.T. -0.029 min  
 Lab File: 3g12912.D  
 Acq: 10 Jan 13 3:08 pm

Tgt Ion: 276 Resp: 805  
 Ion Ratio Lower Upper  
 276 100  
 138 39.8 15.1 55.1  
 277 14.7 3.3 43.3  
 274 46.7 1.5 41.5#



Abundance



**Manual Integrations  
APPROVED  
(compounds with "m" flag)**  
**Judy Nelson  
01/11/13 11:06**

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\011013\  
Data File : 3g12922.D  
Acq On : 10 Jan 2013 7:08 pm  
Operator : DONC  
Sample : D42316-1  
Misc : OP7200,E3G618,30.03,,,1,1  
ALS Vial : 14 Sample Multiplier: 1

Quant Time: Jan 11 08:43:25 2013  
Quant Method : C:\msdchem\1\METHODS\SIMPE3G611.M  
Quant Title : PAHSIM BASE  
QLast Update : Thu Jan 10 14:18:35 2013  
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Naphthalene-d8	5.621	136	62972	4.0000	ug/mL	0.00
6) Acenaphthene-d10	7.349	164	97434m	4.0000	ug/mL	0.02
15) Phenanthrene-d10	8.828	188	152006	4.0000	ug/mL	0.02
19) Chrysene-d12	11.457	240	93621	4.0000	ug/mL	0.01
24) Perylene-d12	12.820	264	94494	4.0000	ug/mL	0.01

System Monitoring Compounds

2) Nitrobenzene-d5	4.935	82	238176	42.0496	ug/mL	-0.01
Spiked Amount	50.000	Range	25 - 135	Recovery	=	84.10%
7) 2-Fluorobiphenyl	6.676	172	979998	25.0126	ug/mL	0.00
Spiked Amount	50.000	Range	25 - 135	Recovery	=	50.02%
21) Terphenyl-d14	10.418	244	462700	36.3216	ug/mL	0.02
Spiked Amount	50.000	Range	25 - 135	Recovery	=	72.64%

Target Compounds Qvalue

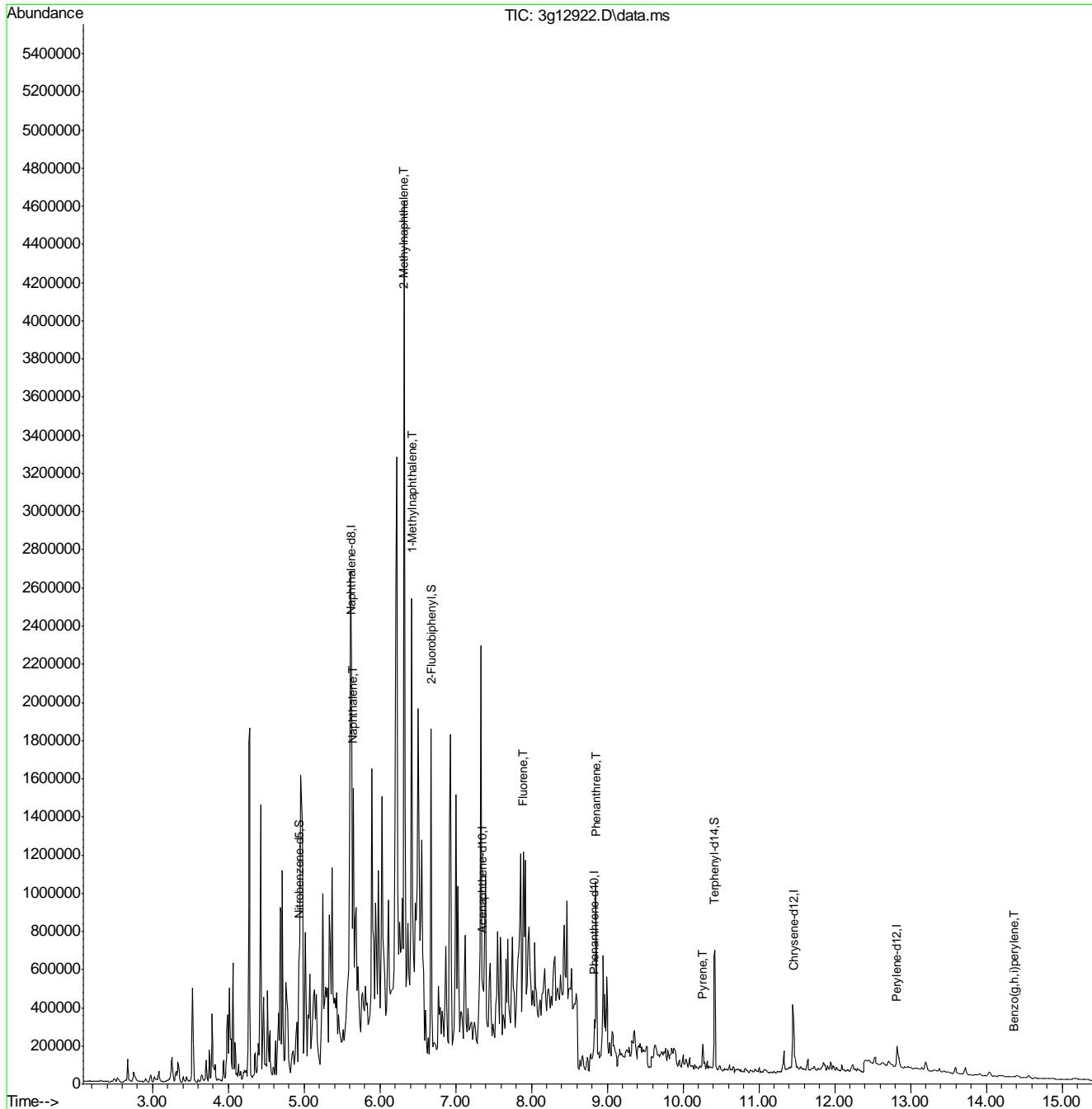
3) N-Nitrosodimethylamine	2.385	74	102	N.D.		
4) N-Nitrosodi-propylamine	0.000	70	0	N.D.	d	
5) Naphthalene	5.646	128	911543	50.4507	ug/mL#	57
8) 2-Methylnaphthalene	6.319	142	1841396	58.8127	ug/mL	95
9) 1-Methylnaphthalene	6.419	142	786800m	28.7316	ug/mL	
10) Acenaphthylene	0.000	152	0	N.D.	d	
11) Acenaphthene	0.000	154	0	N.D.	d	
12) Dibenzofuran	0.000	168	0	N.D.	d	
13) Fluorene	7.893	166	325715m	8.3108	ug/mL	
14) Diphenylamine	0.000	169	0	N.D.	d	
16) Phenanthrene	8.851	178	623923	10.6215	ug/mL	80
17) Anthracene	0.000	178	0	N.D.	d	
18) Fluoranthene	0.000	202	0	N.D.	d	
20) Pyrene	10.260	202	76001	1.5201	ug/mL#	57
22) Benzo(a)anthracene	0.000	228	0	N.D.	d	
23) Chrysene	0.000	228	0	N.D.	d	
25) Benzo(b)fluoranthene	0.000	252	0	N.D.	d	
26) Benzo(k)fluoranthene	0.000	252	0	N.D.	d	
27) Benzo(a)pyrene	0.000	252	0	N.D.	d	
28) Indeno(1,2,3-cd)pyrene	0.000	276	0	N.D.	d	
29) Dibenz(a,h)anthracene	0.000	278	0	N.D.	d	
30) Benzo(g,h,i)perylene	14.366	276	4649	0.1067	ug/mL#	80

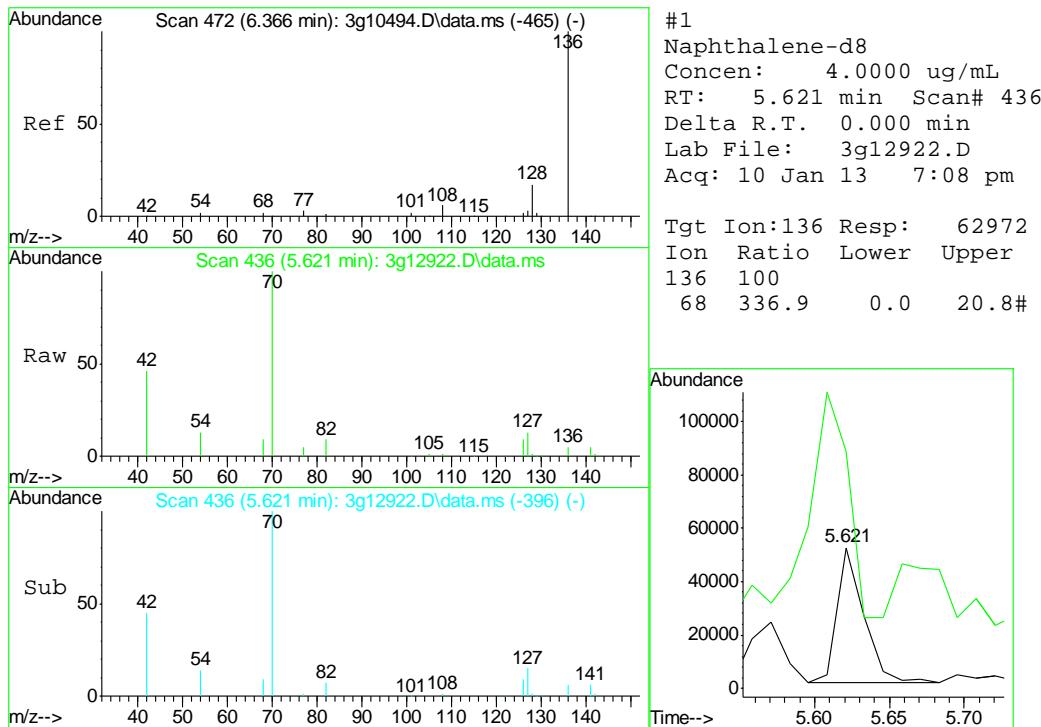
(#) = qualifier out of range (m) = manual integration (+) = signals summed

## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\011013\  
 Data File : 3g12922.D  
 Acq On : 10 Jan 2013 7:08 pm  
 Operator : DONC  
 Sample : D42316-1  
 Misc : OP7200,E3G618,30.03,,,1,1  
 ALS Vial : 14 Sample Multiplier: 1

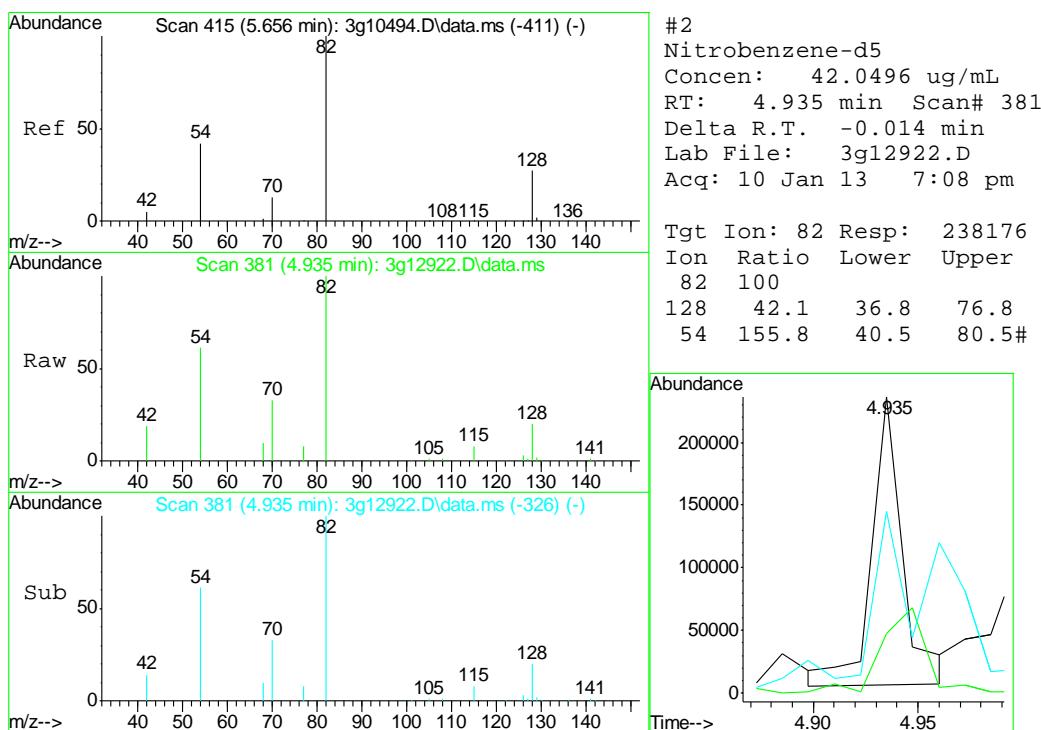
Quant Time: Jan 11 08:43:25 2013  
 Quant Method : C:\msdchem\1\METHODS\SIMPE3G611.M  
 Quant Title : PAHSIM BASE  
 QLast Update : Thu Jan 10 14:18:35 2013  
 Response via : Initial Calibration

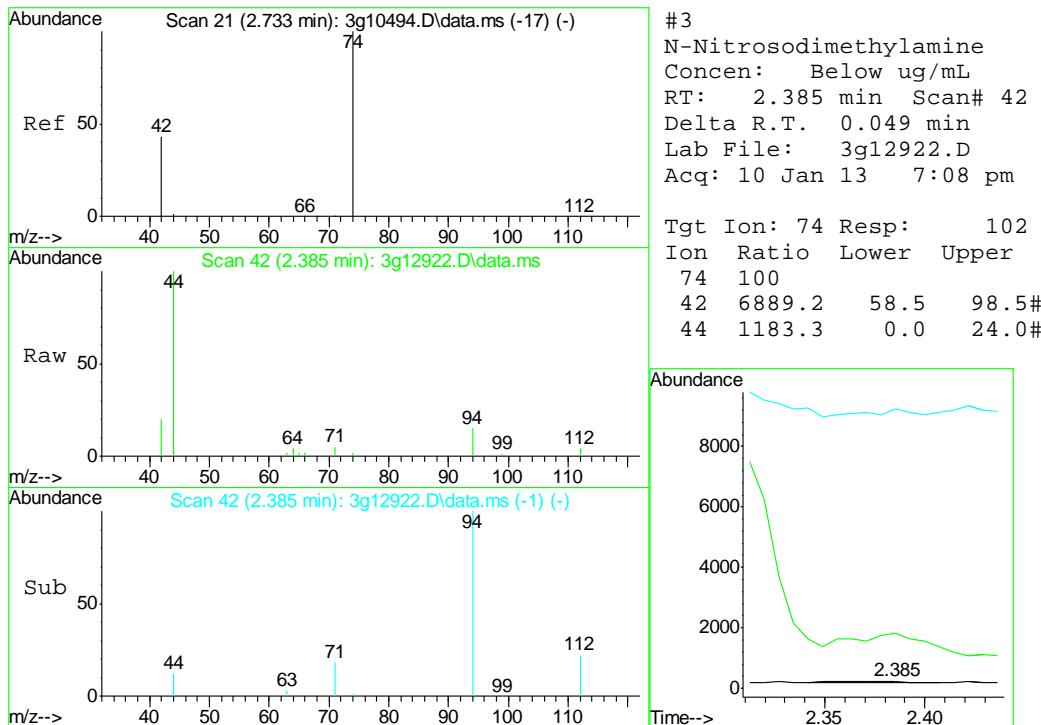




9.1.2

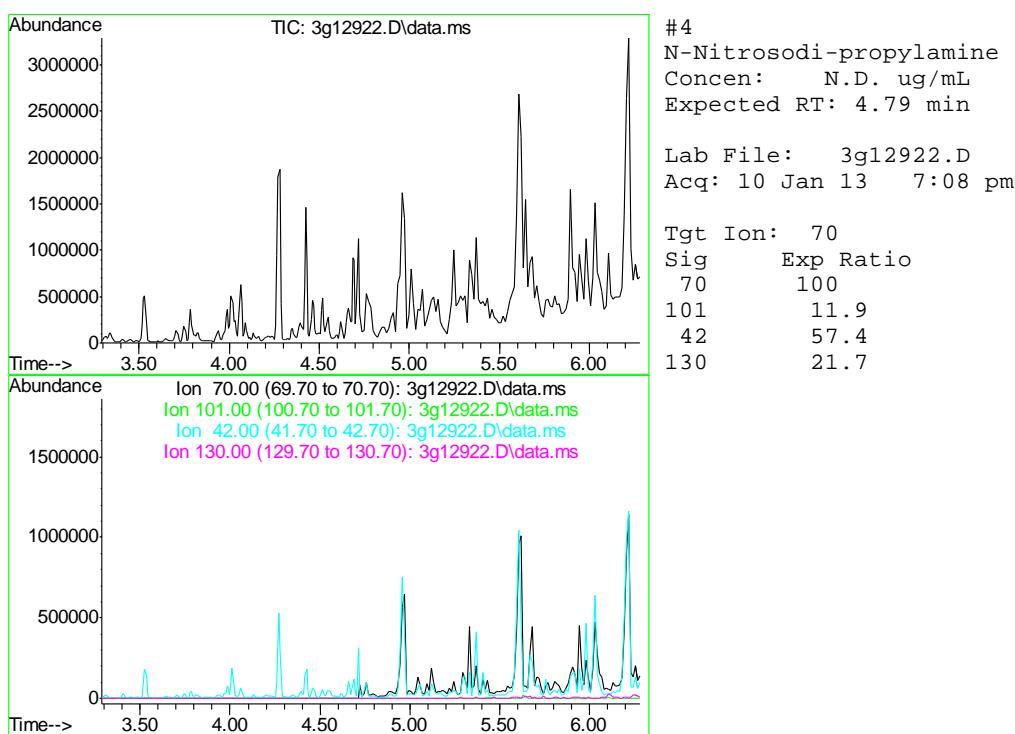
9

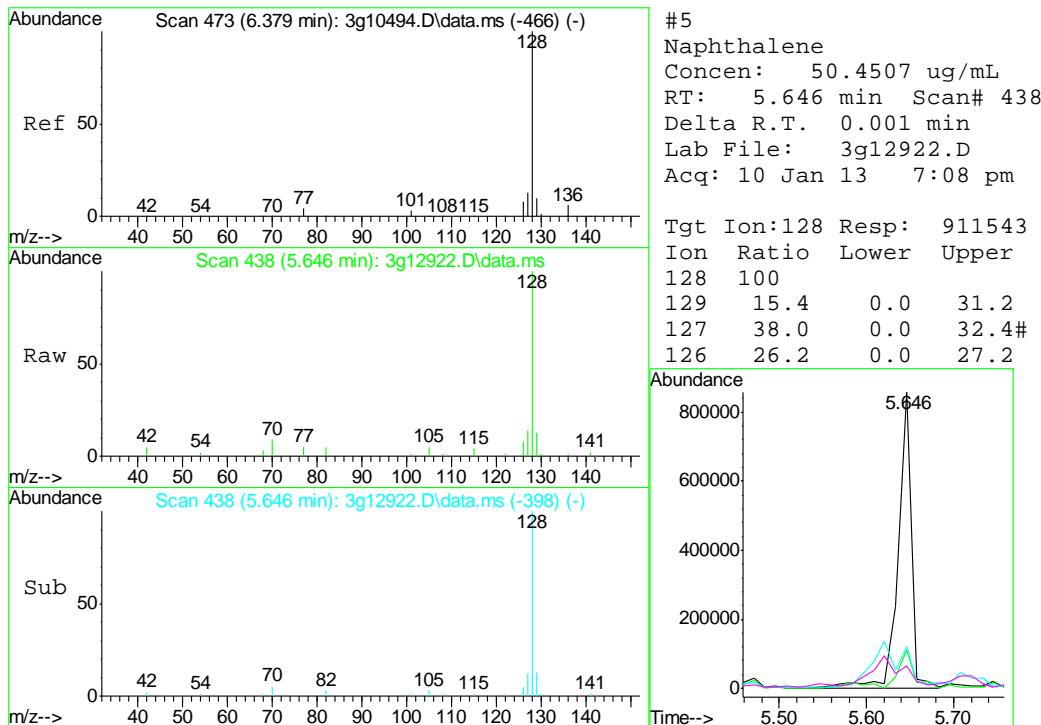




9.1.2

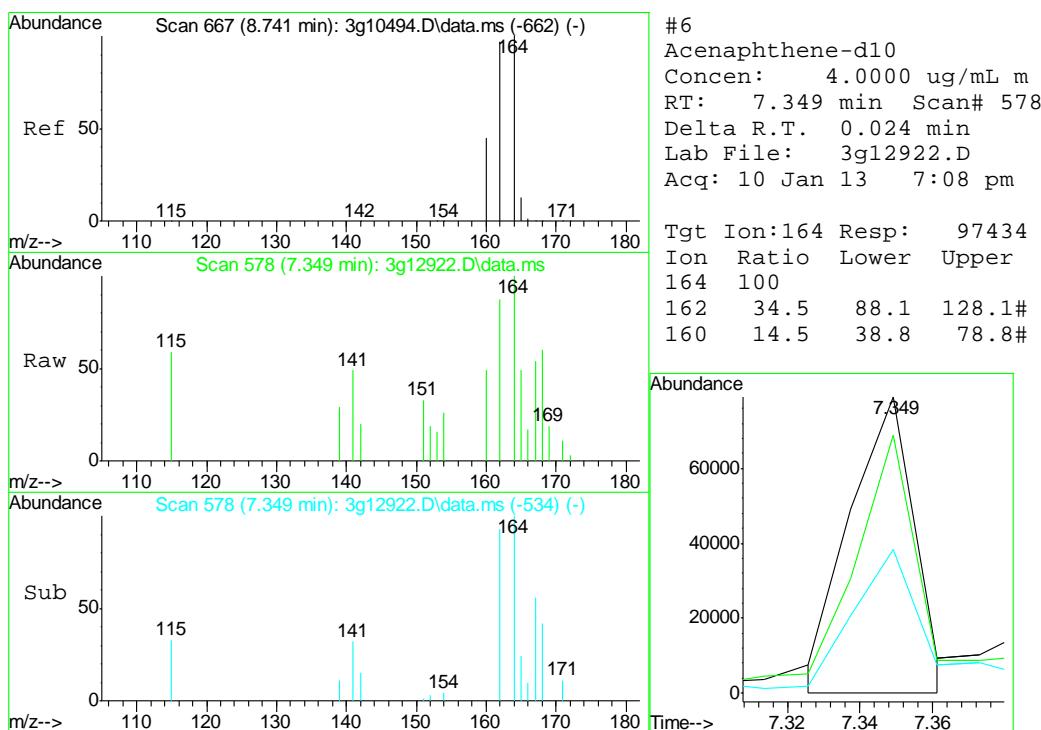
9

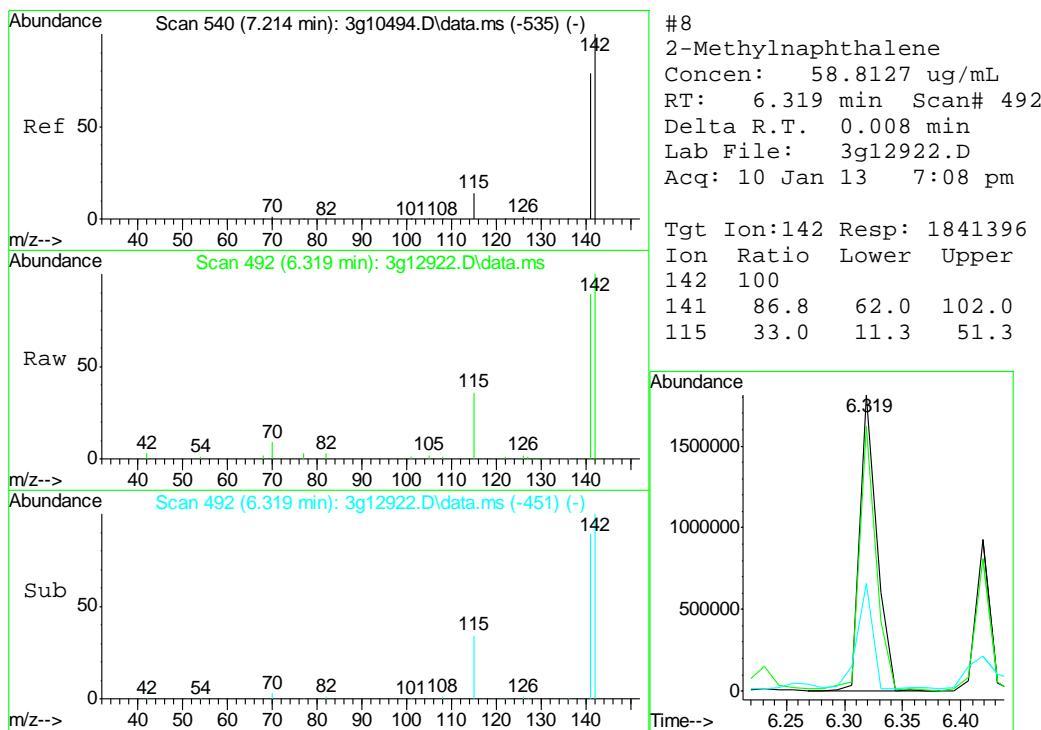
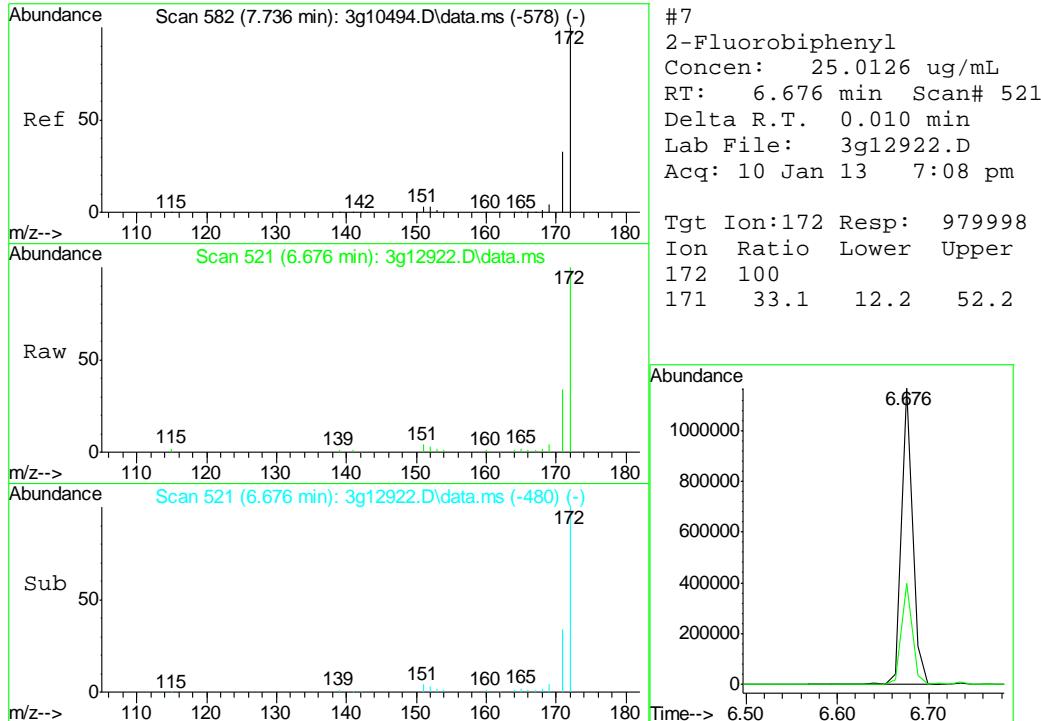


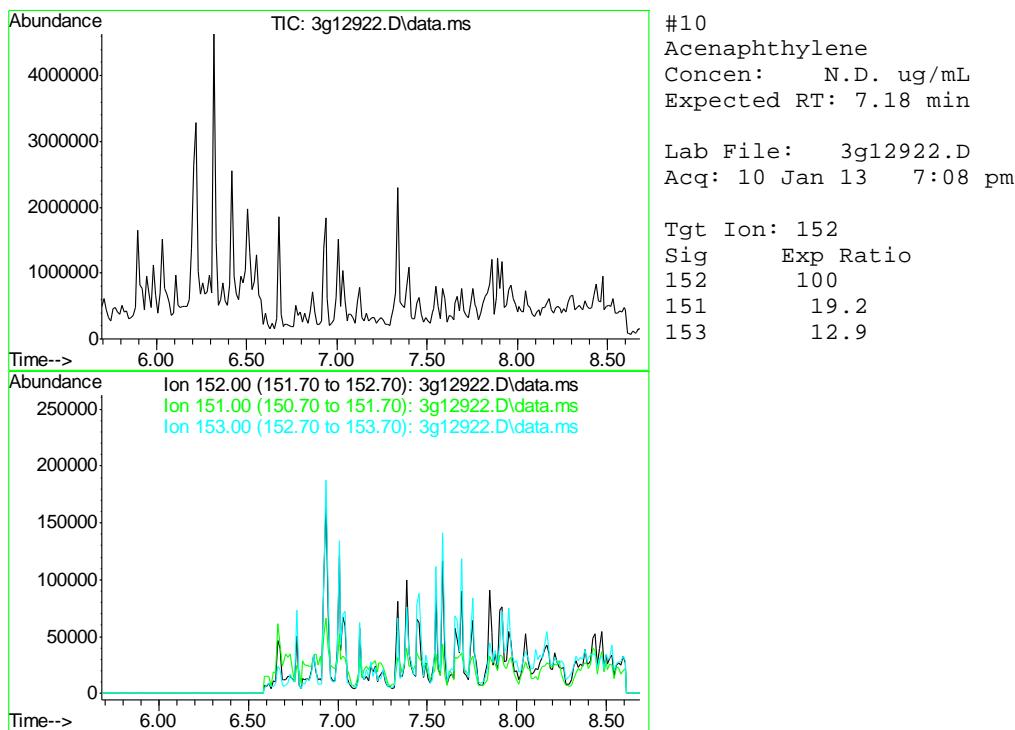
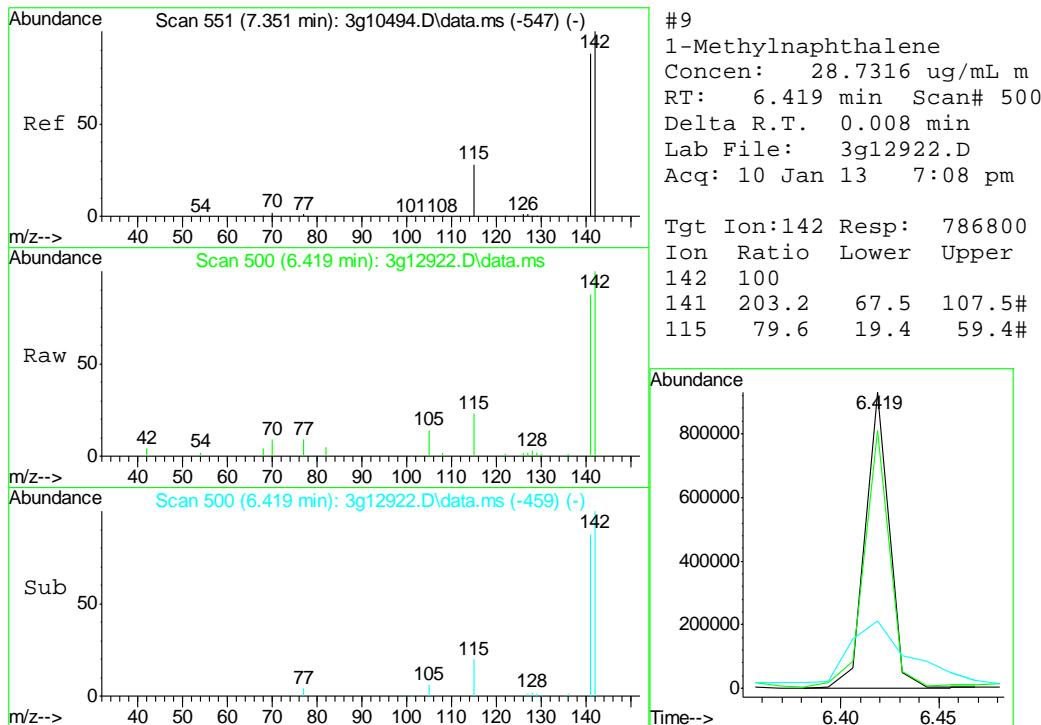


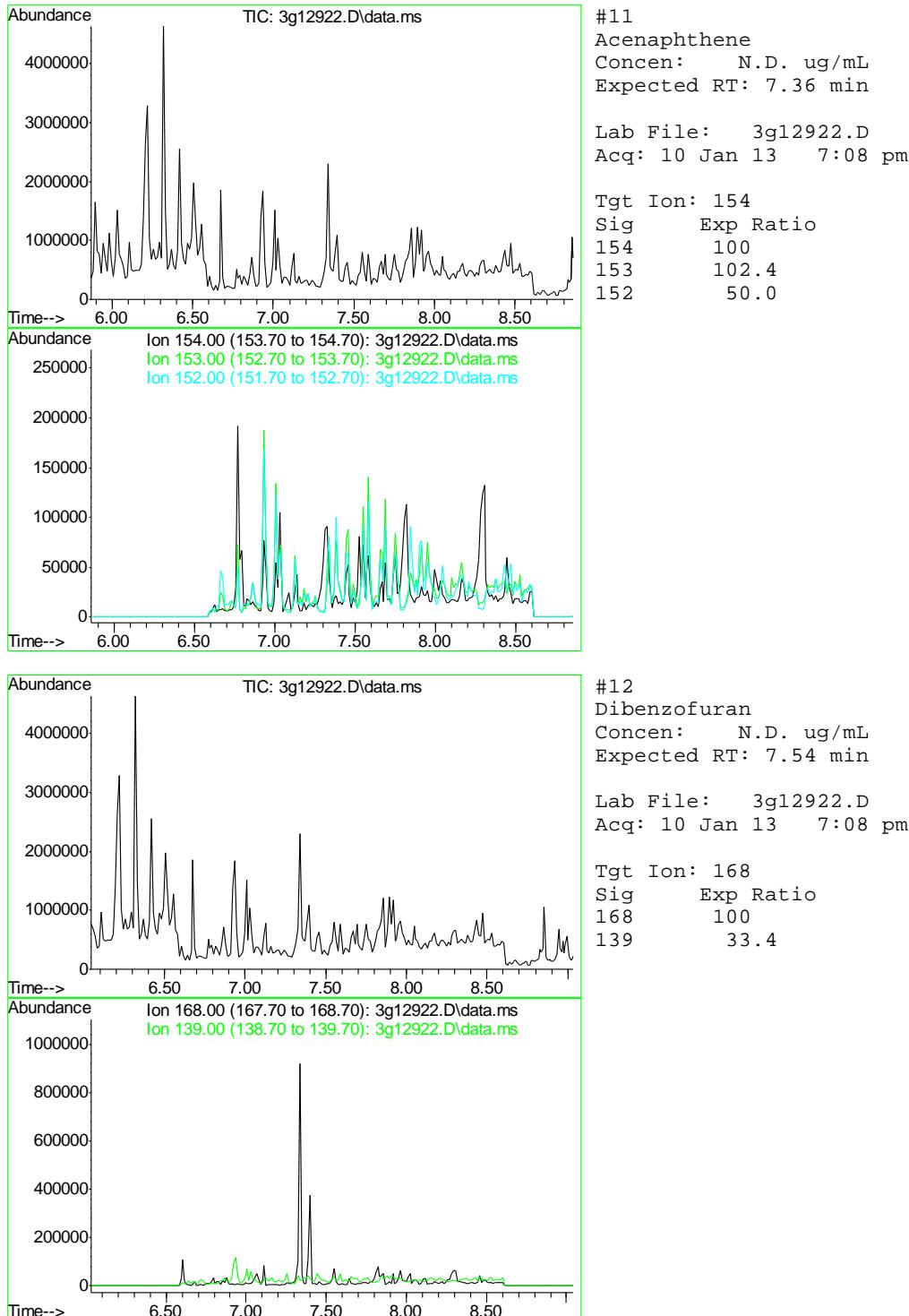
9.1.2

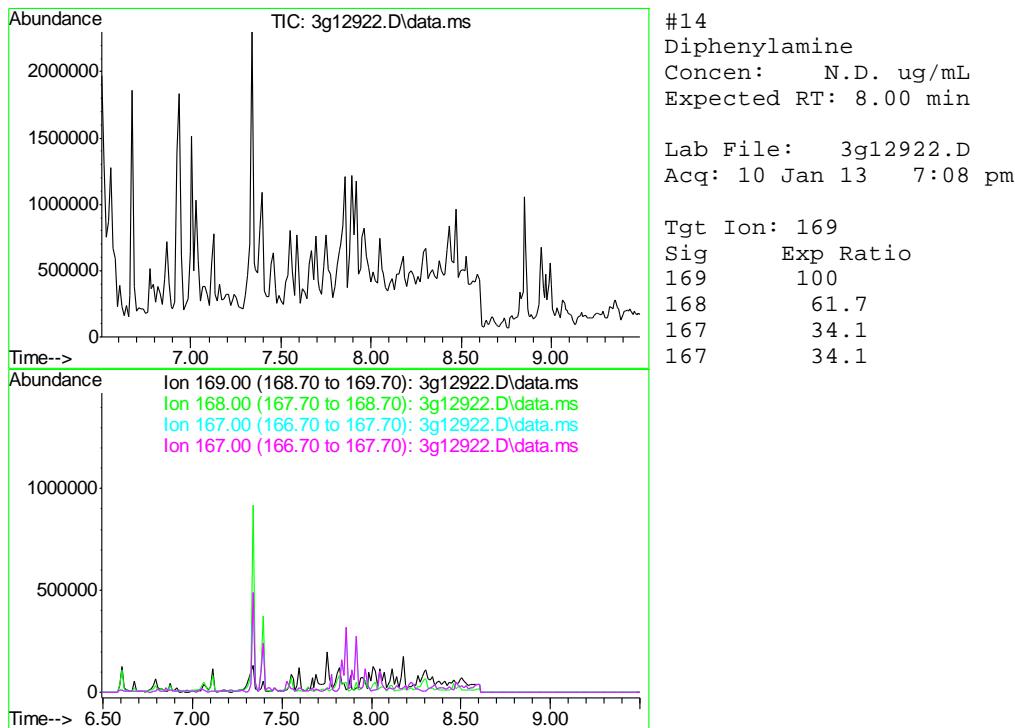
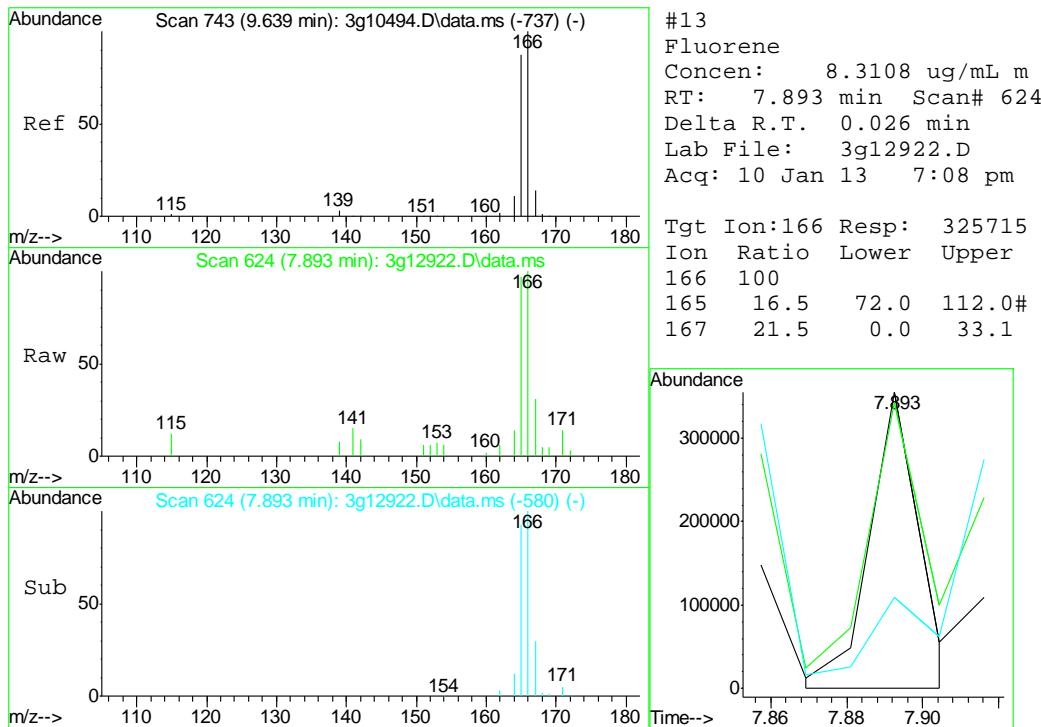
9

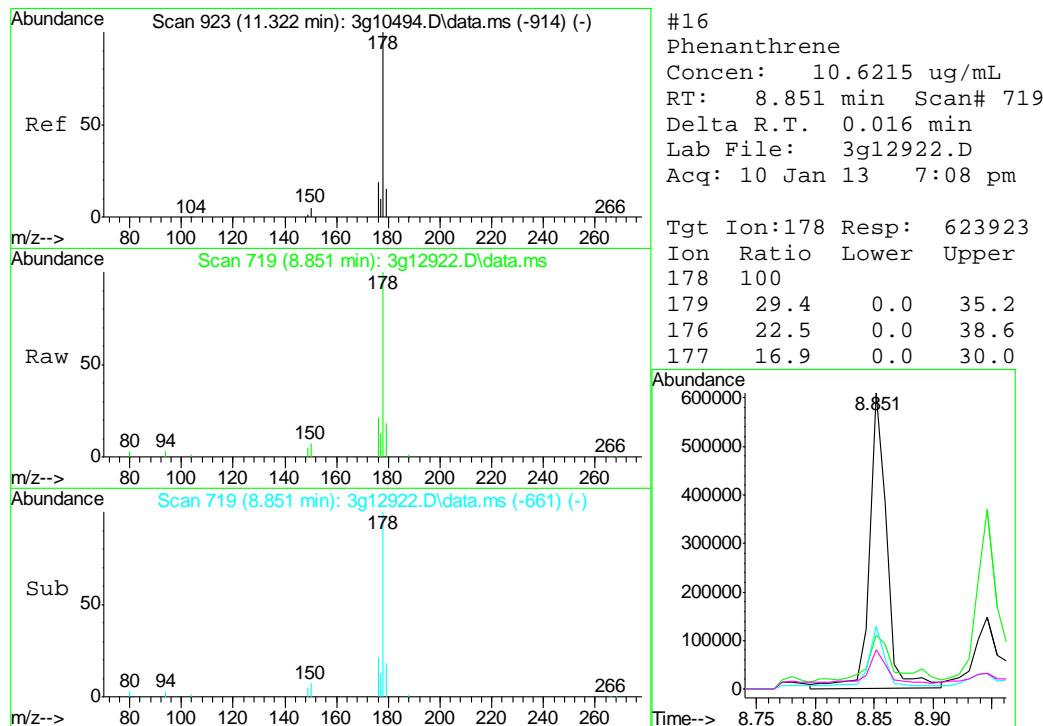
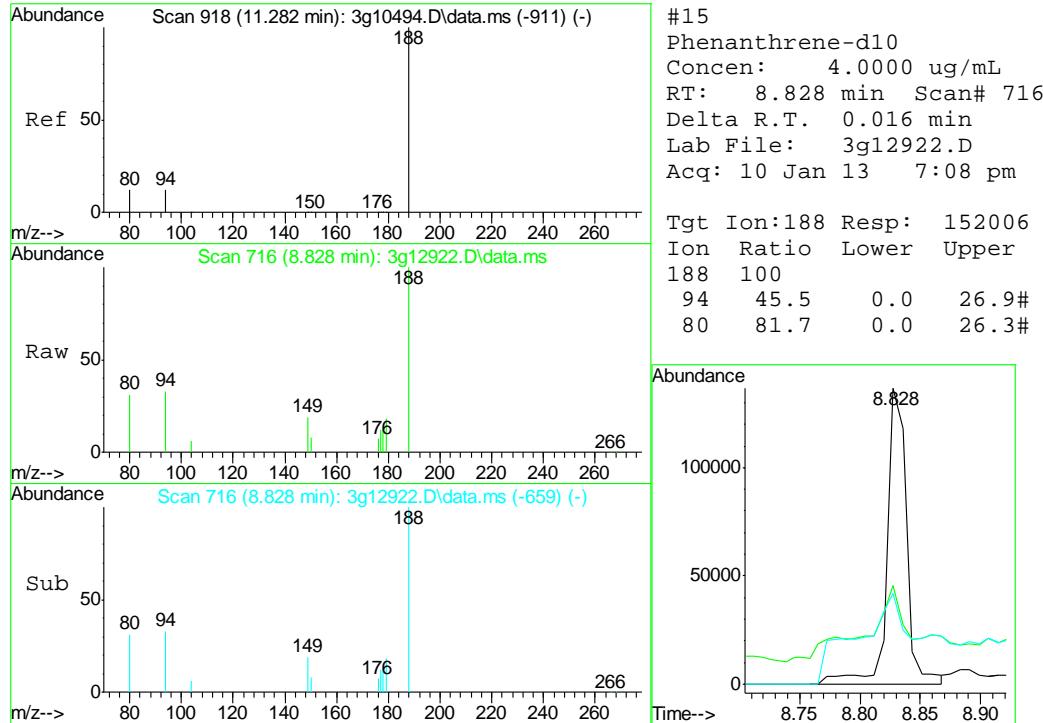


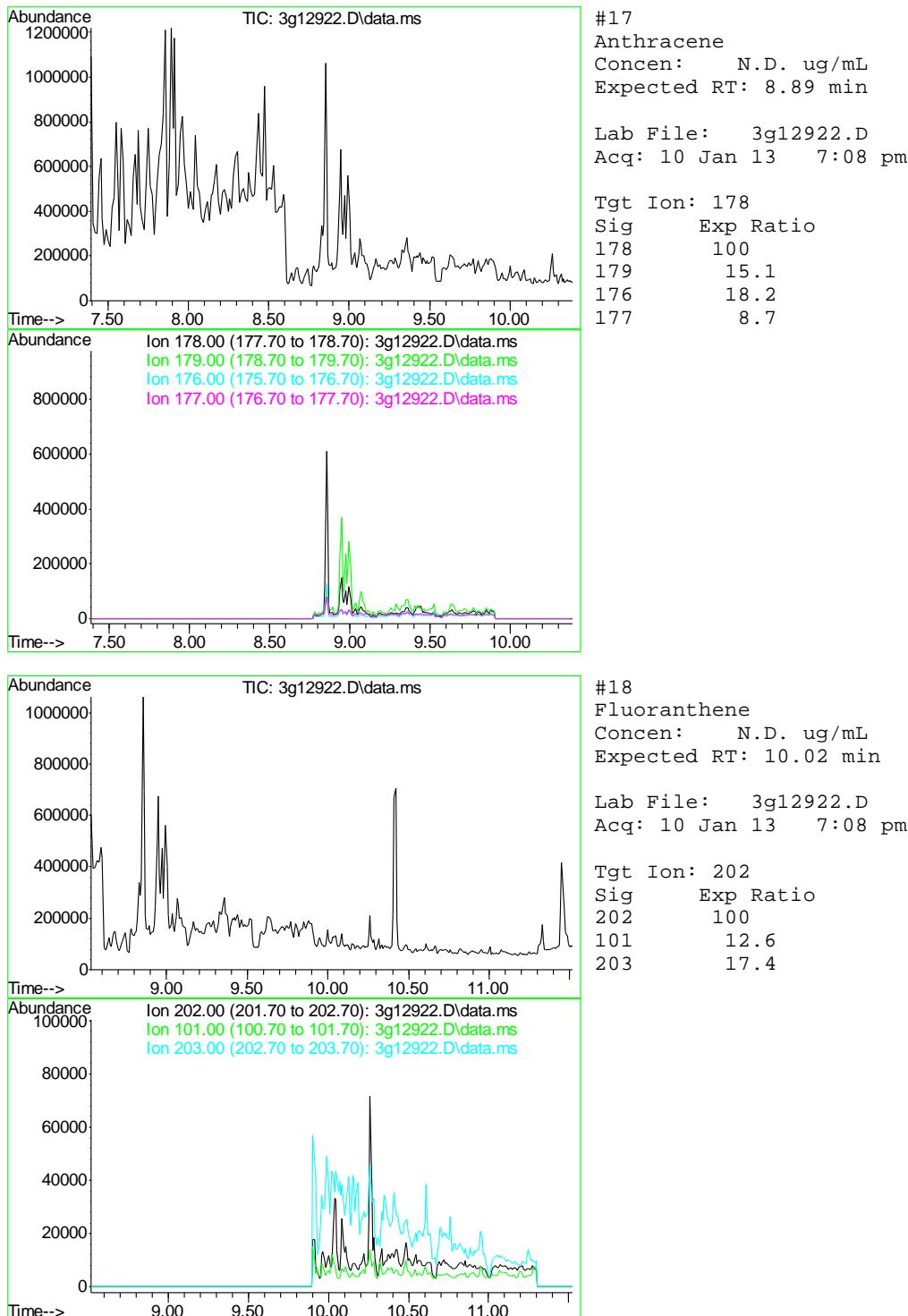


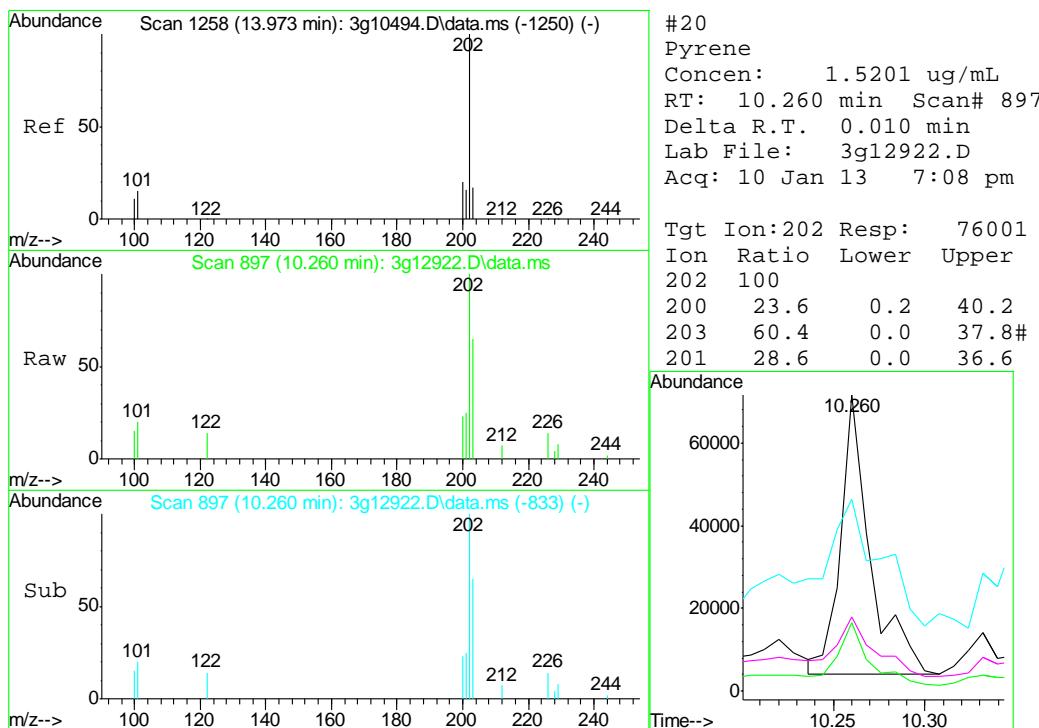
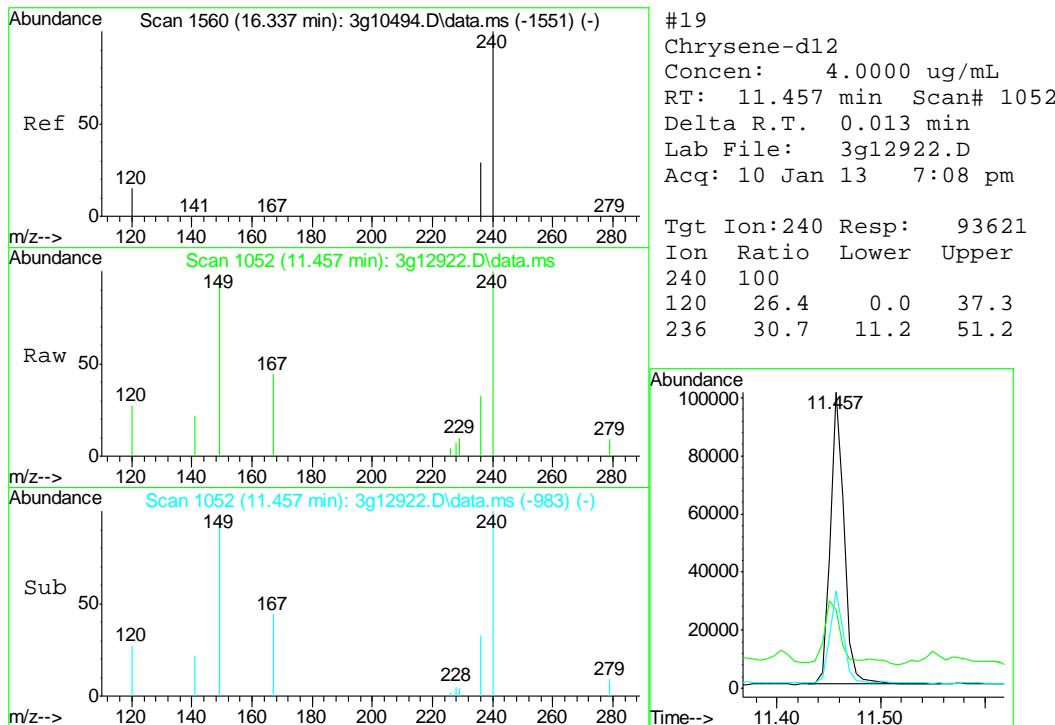


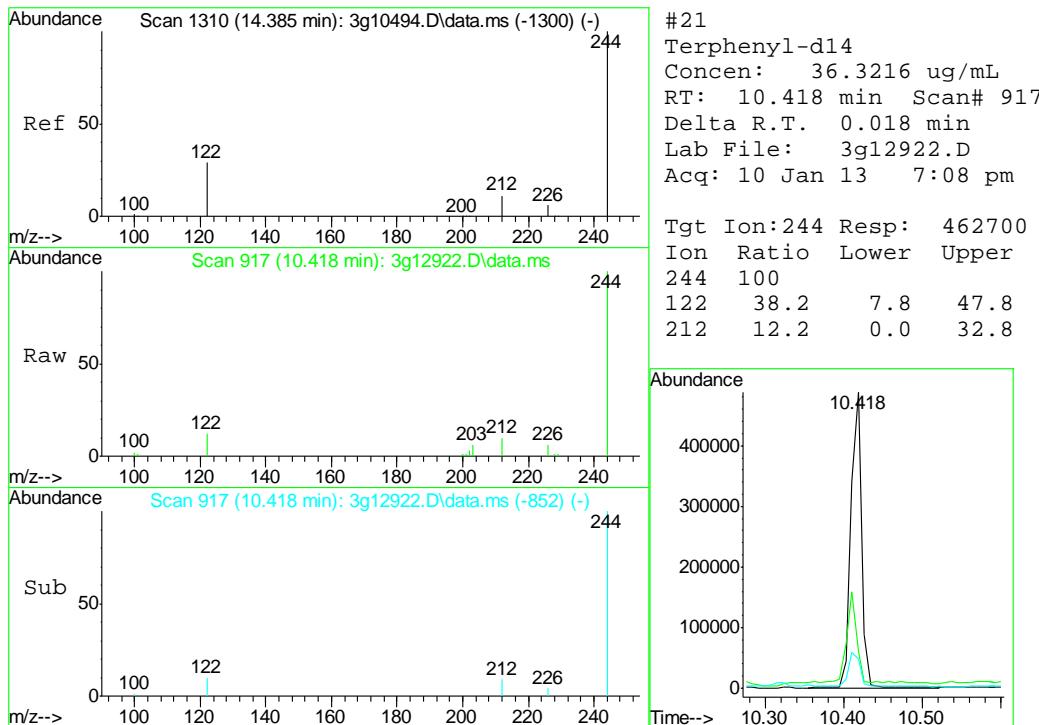
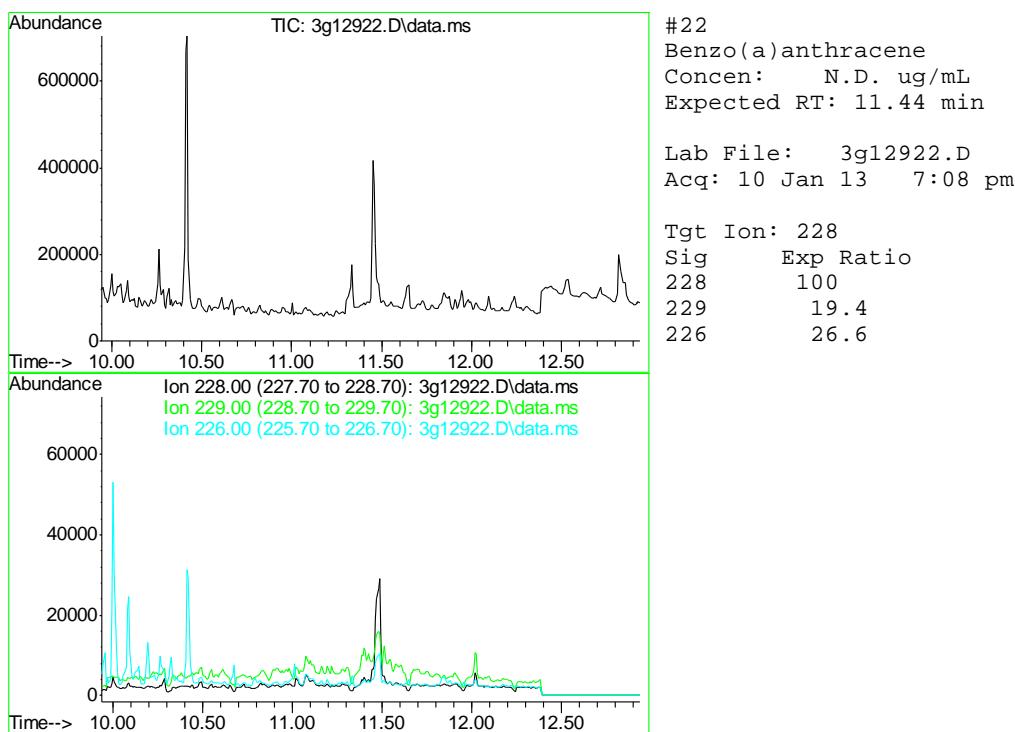


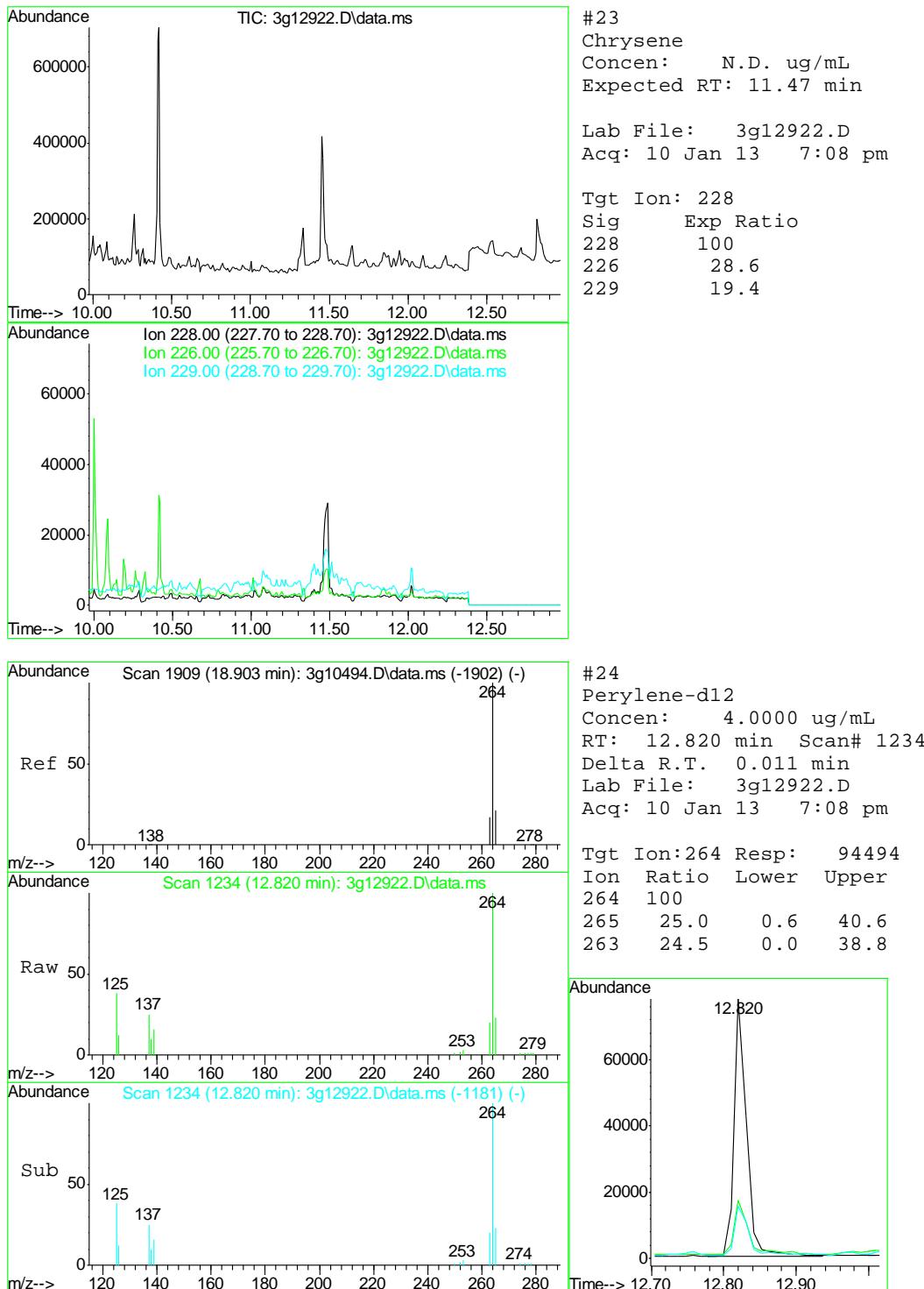


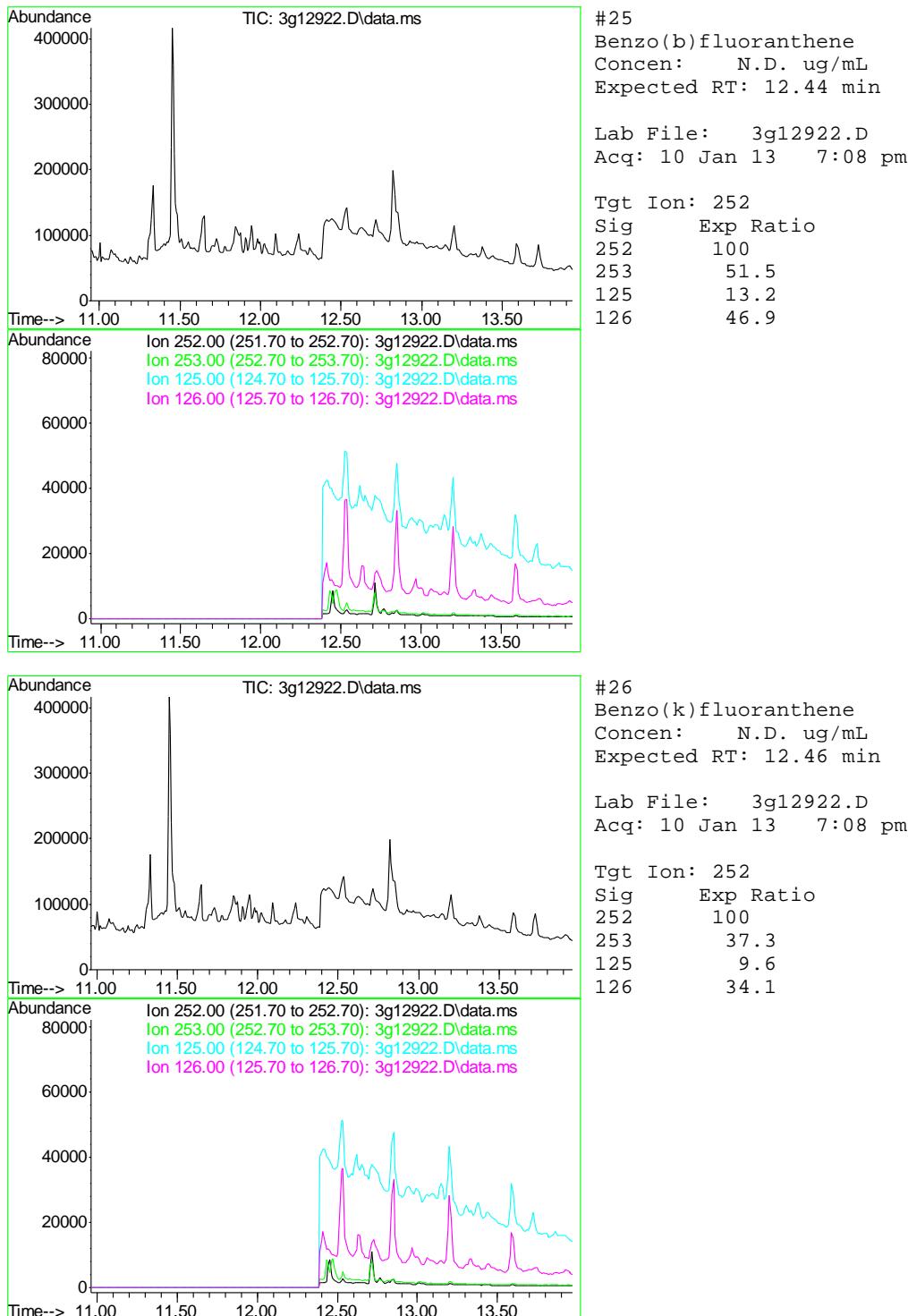


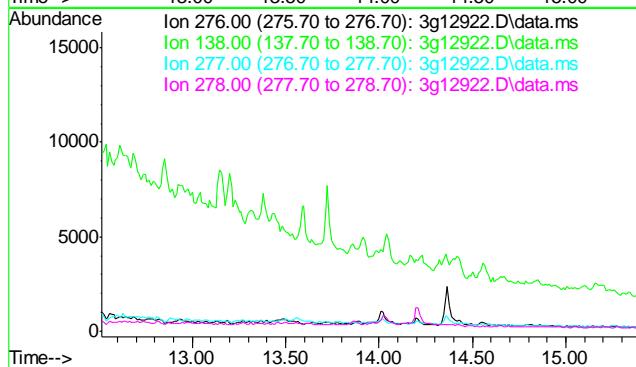
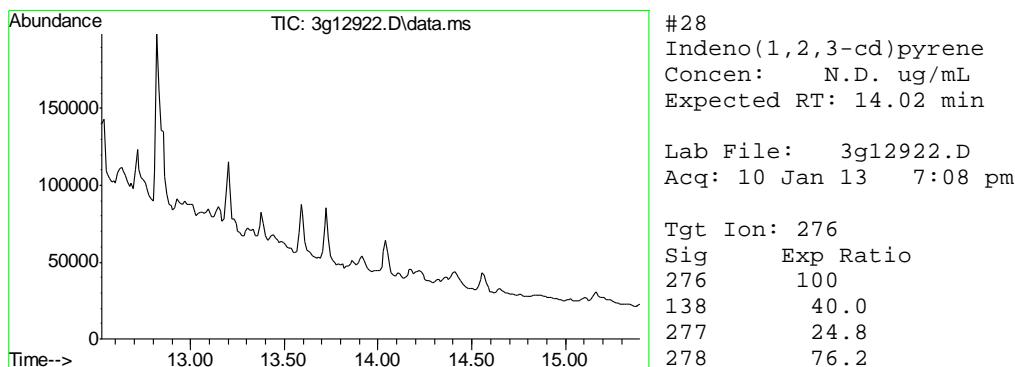
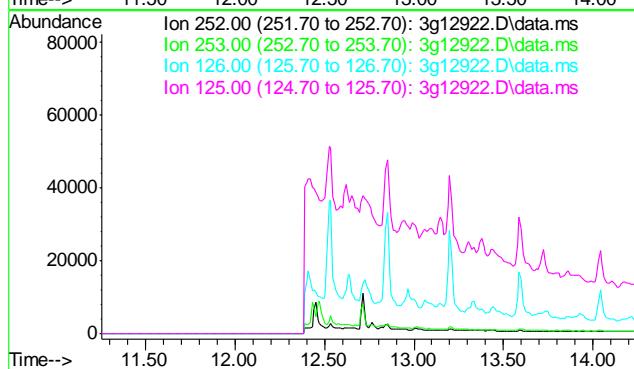
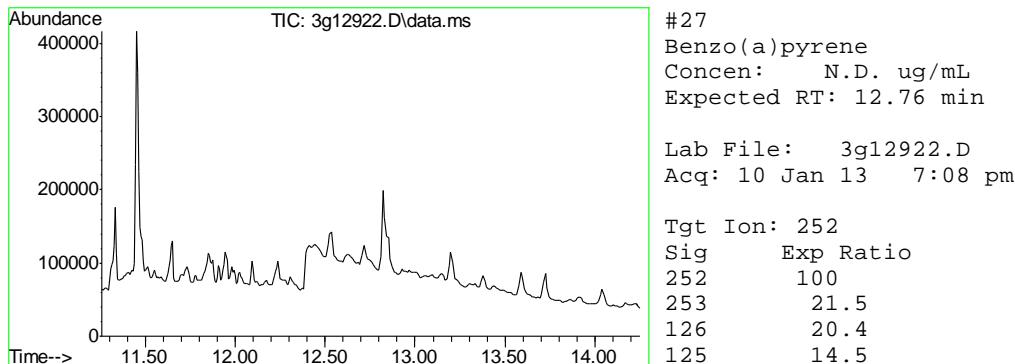


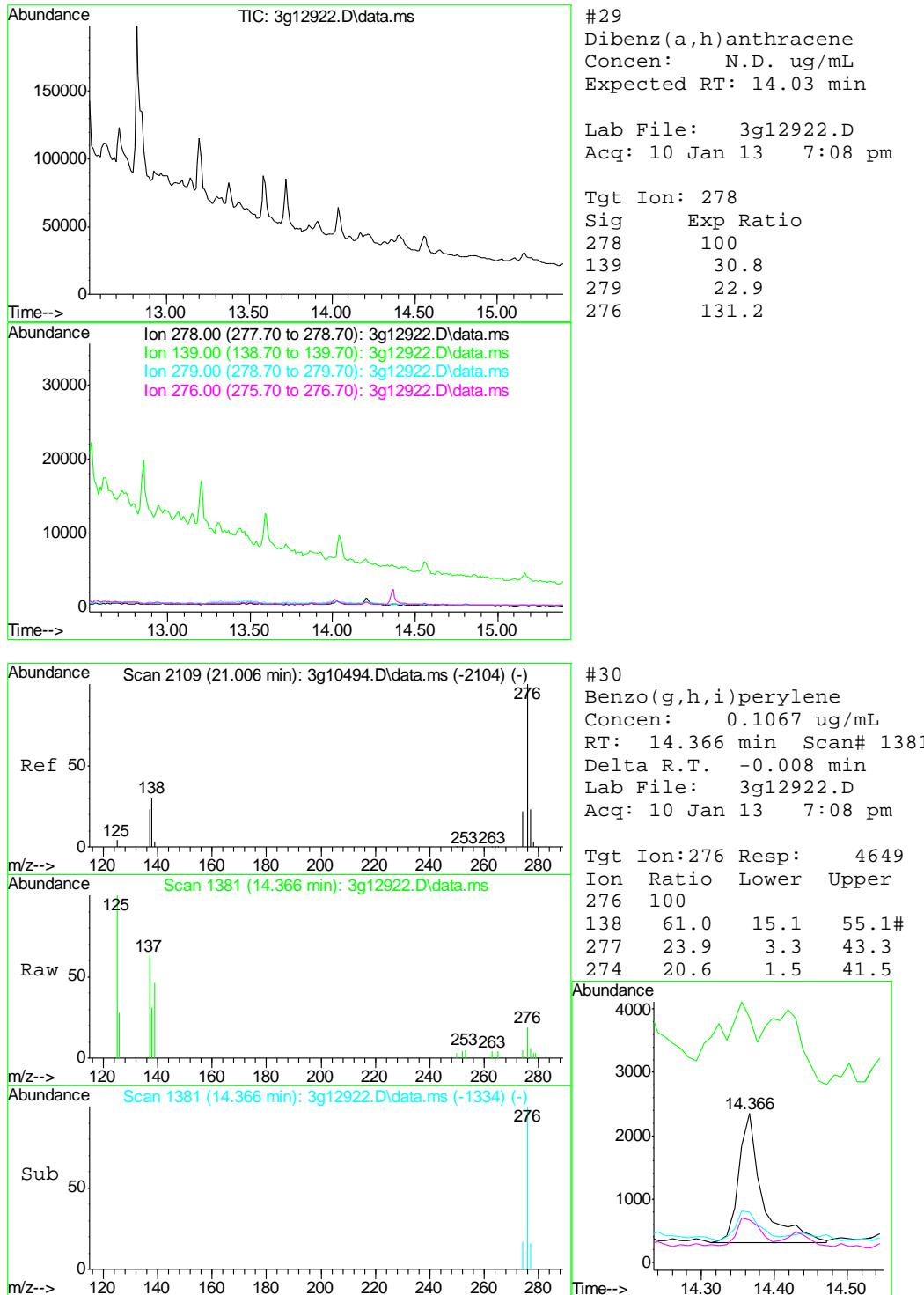


9.1.2  
9









## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\011013\  
 Data File : 3g12911.D  
 Acq On : 10 Jan 2013 2:44 pm  
 Operator : DONC  
 Sample : OP7200-MB  
 Misc : OP7200,E3G618,30.00,,,1,1  
 ALS Vial : 5 Sample Multiplier: 1

Quant Time: Jan 10 15:58:55 2013  
 Quant Method : C:\msdchem\1\METHODS\SIMPE3G611.M  
 Quant Title : PAHSIM BASE  
 QLast Update : Thu Jan 10 14:18:35 2013  
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Naphthalene-d8	5.621	136	104180	4.0000	ug/mL	0.00
6) Acenaphthene-d10	7.326	164	59648	4.0000	ug/mL	0.00
15) Phenanthrene-d10	8.812	188	94038	4.0000	ug/mL	0.00
19) Chrysene-d12	11.444	240	52962	4.0000	ug/mL	0.00
24) Perylene-d12	12.810	264	32109	4.0000	ug/mL	0.00

## System Monitoring Compounds

2) Nitrobenzene-d5	4.935	82	491377	52.4375	ug/mL	-0.01
Spiked Amount	50.000	Range	25 - 135	Recovery	=	104.88%
7) 2-Fluorobiphenyl	6.664	172	1123238	50.4729	ug/mL	0.00
Spiked Amount	50.000	Range	25 - 135	Recovery	=	100.94%
21) Terphenyl-d14	10.403	244	575142	79.8087	ug/mL	0.00
Spiked Amount	50.000	Range	25 - 135	Recovery	=	159.62%#

## Target Compounds

				Qvalue
3) N-Nitrosodimethylamine	2.291	74	47	N.D.
4) N-Nitrosodi-propylamine	0.000	70	0	N.D. d
5) Naphthalene	5.633	128	455	N.D.
8) 2-Methylnaphthalene	6.306	142	230	N.D.
9) 1-Methylnaphthalene	6.406	142	73	N.D.
10) Acenaphthylene	7.385	152	128	N.D.
11) Acenaphthene	7.326	154	304	N.D.
12) Dibenzofuran	7.538	168	61	N.D.
13) Fluorene	0.000	166	0	N.D. d
14) Diphenylamine	0.000	169	0	N.D. d
16) Phenanthrene	8.835	178	183	N.D.
17) Anthracene	8.875	178	102	N.D.
18) Fluoranthene	9.928	202	200	N.D.
20) Pyrene	0.000	202	0	N.D. d
22) Benzo(a)anthracene	11.444	228	289	N.D.
23) Chrysene	11.444	228	289	N.D.
25) Benzo(b)fluoranthene	12.484	252	140	N.D.
26) Benzo(k)fluoranthene	12.484	252	140	N.D.
27) Benzo(a)pyrene	0.000	252	0	N.D. d
28) Indeno(1,2,3-cd)pyrene	14.040	276	37	N.D.
29) Dibenz(a,h)anthracene	0.000	278	0	N.D. d
30) Benzo(g,h,i)perylene	14.345	276	50	N.D.

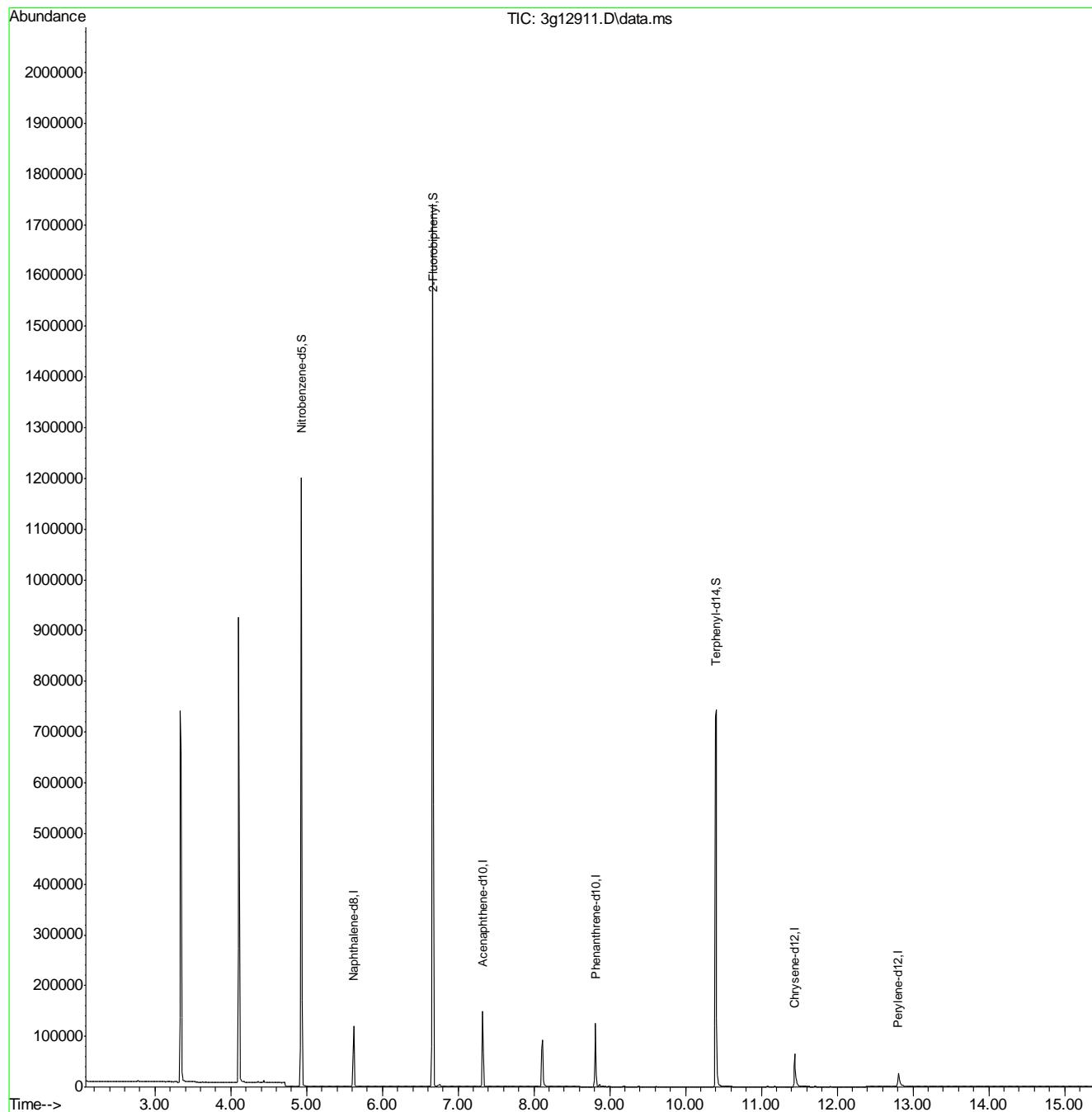
(#) = qualifier out of range (m) = manual integration (+) = signals summed

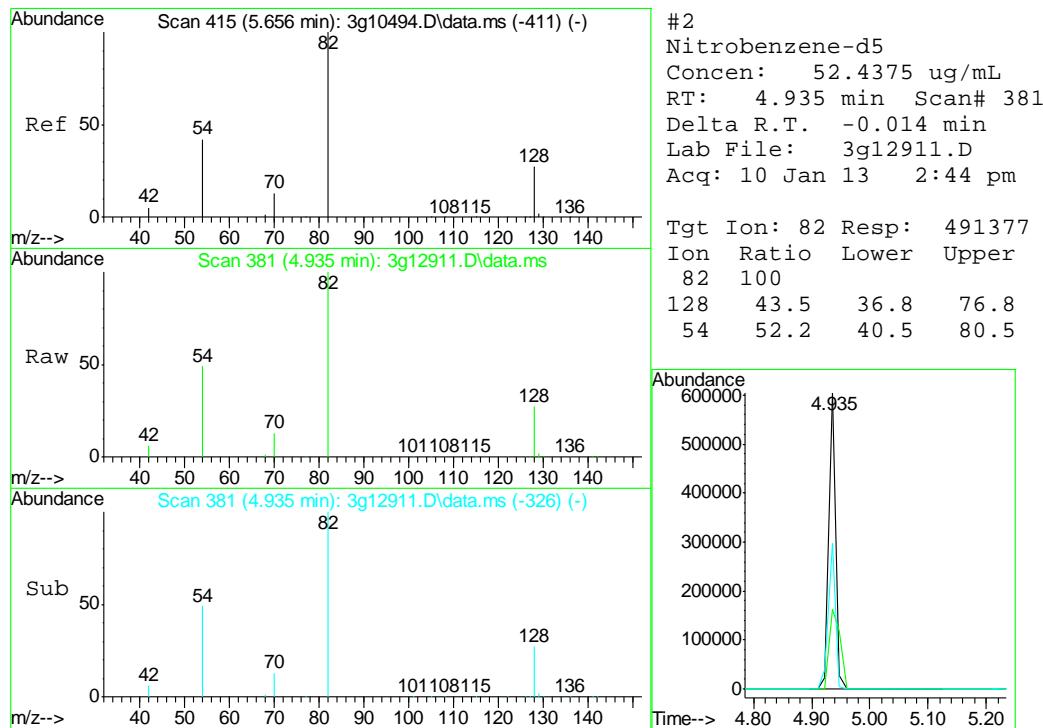
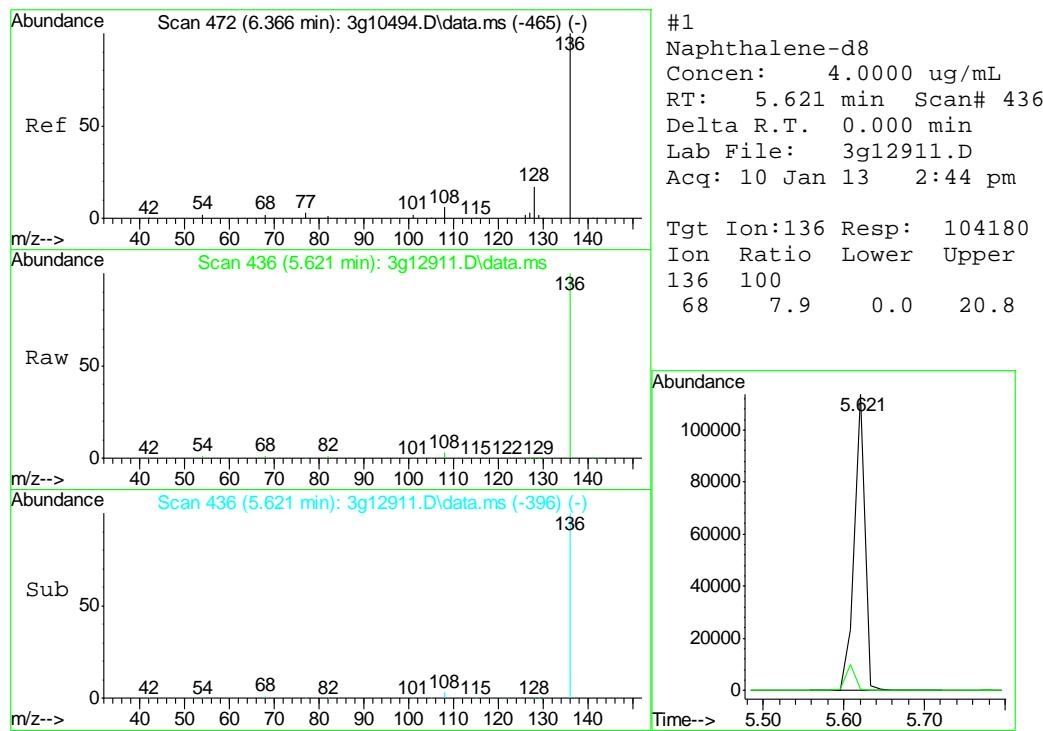
9.2.1  
9

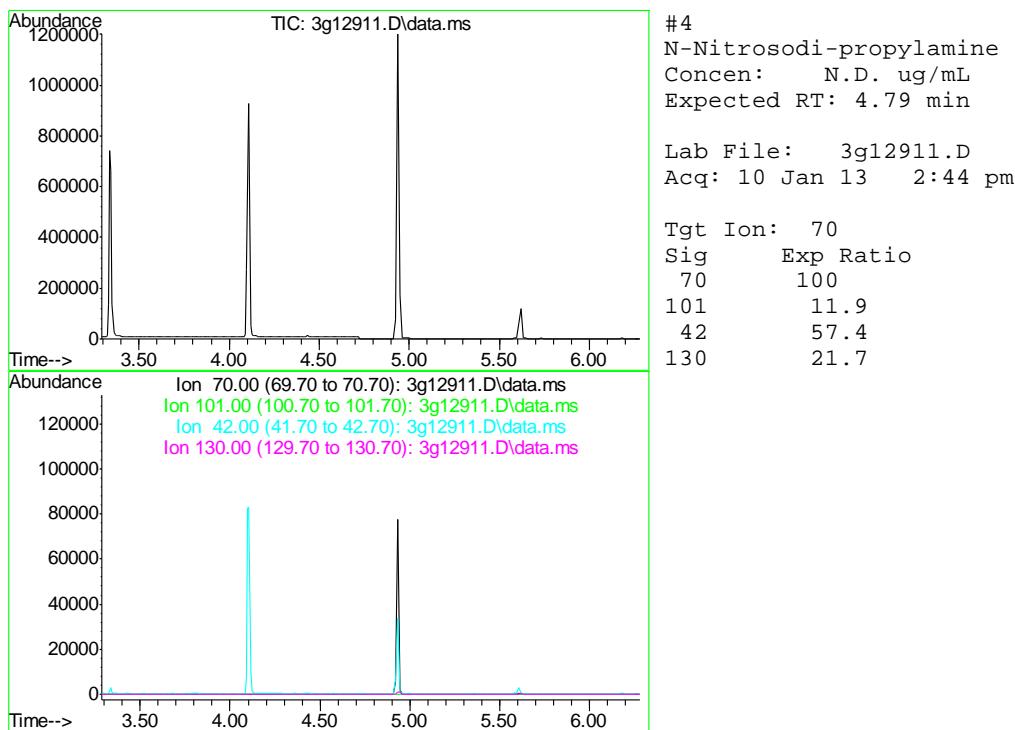
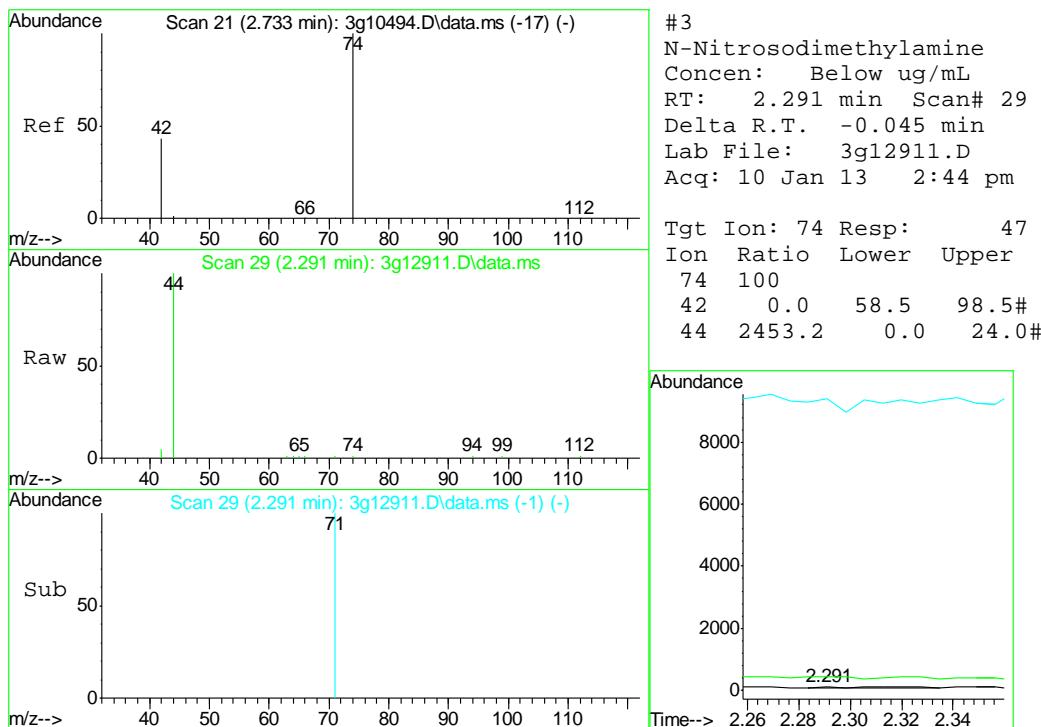
## Quantitation Report (QT Reviewed)

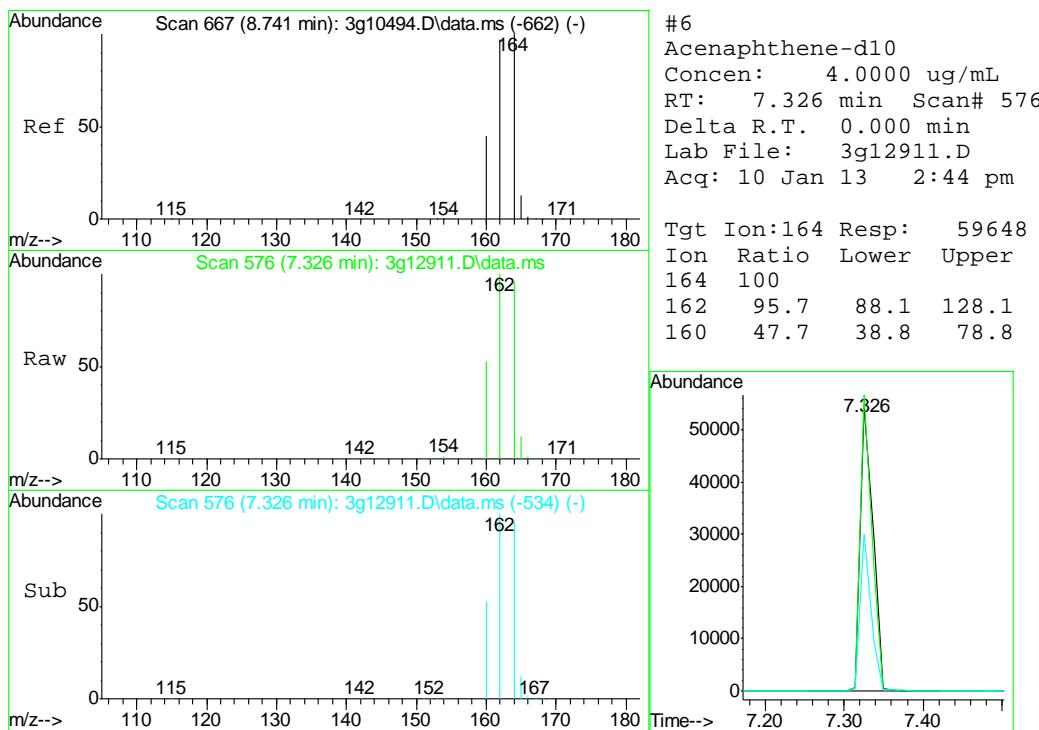
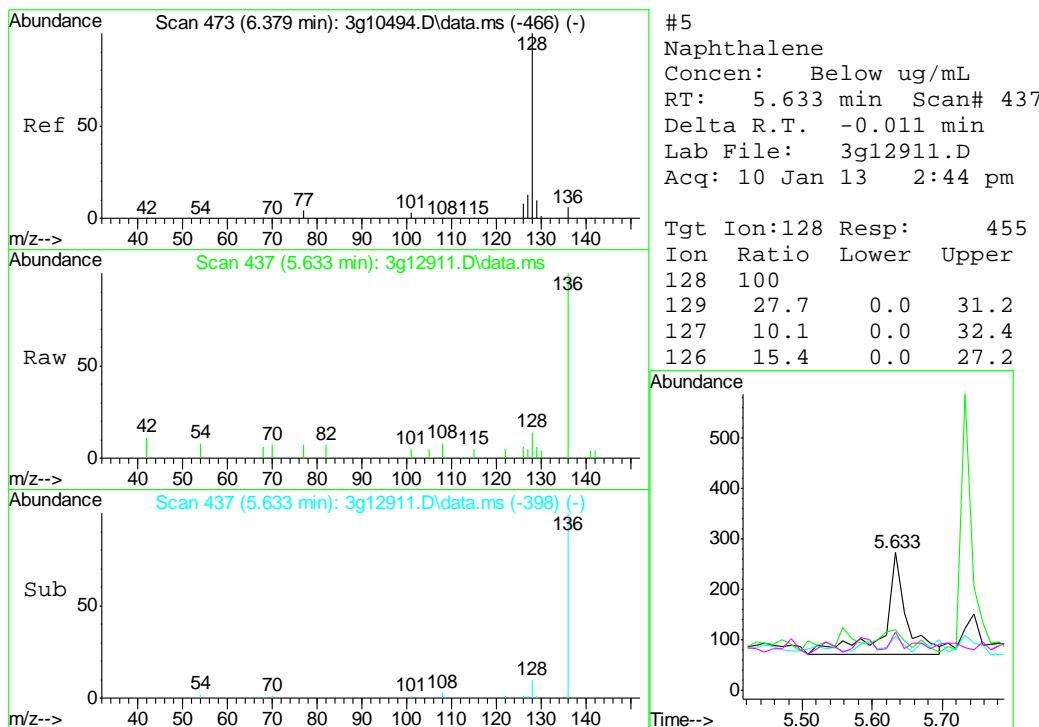
Data Path : C:\msdchem\1\DATA\011013\  
 Data File : 3g12911.D  
 Acq On : 10 Jan 2013 2:44 pm  
 Operator : DONC  
 Sample : OP7200-MB  
 Misc : OP7200,E3G618,30.00,,,1,1  
 ALS Vial : 5 Sample Multiplier: 1

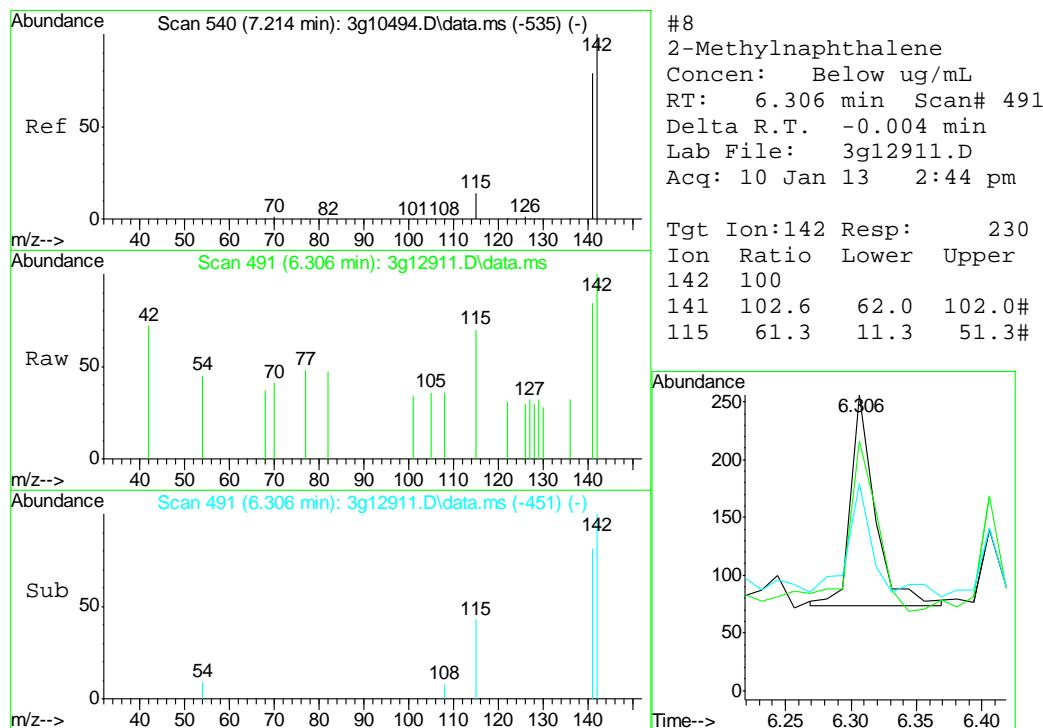
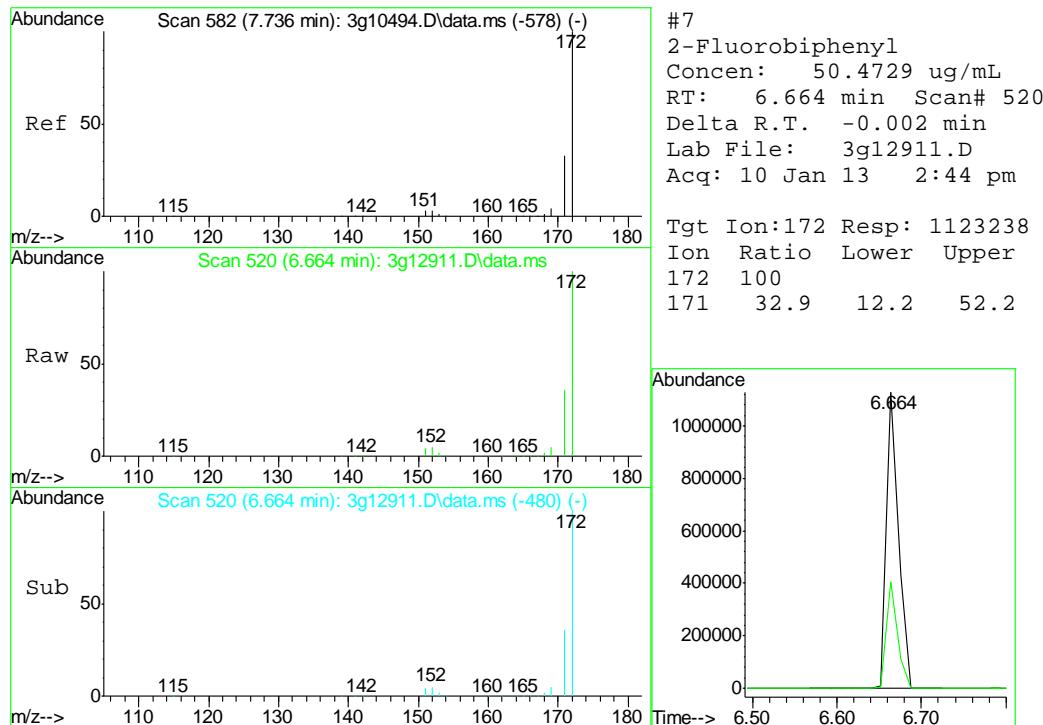
Quant Time: Jan 10 15:58:55 2013  
 Quant Method : C:\msdchem\1\METHODS\SIMPE3G611.M  
 Quant Title : PAHSIM BASE  
 QLast Update : Thu Jan 10 14:18:35 2013  
 Response via : Initial Calibration

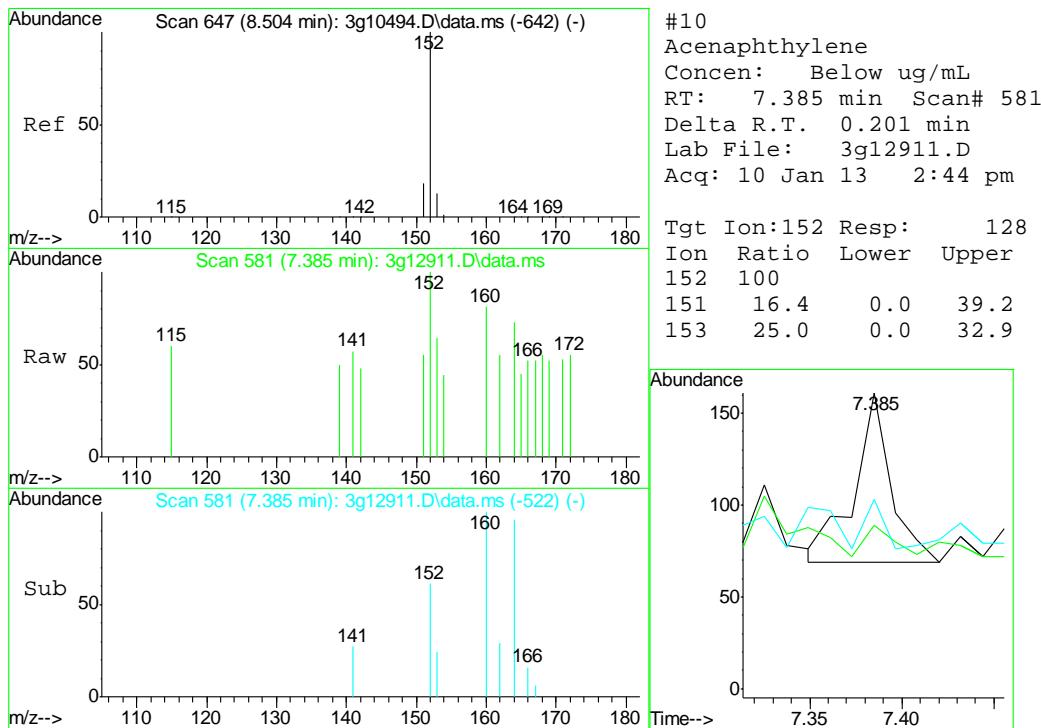
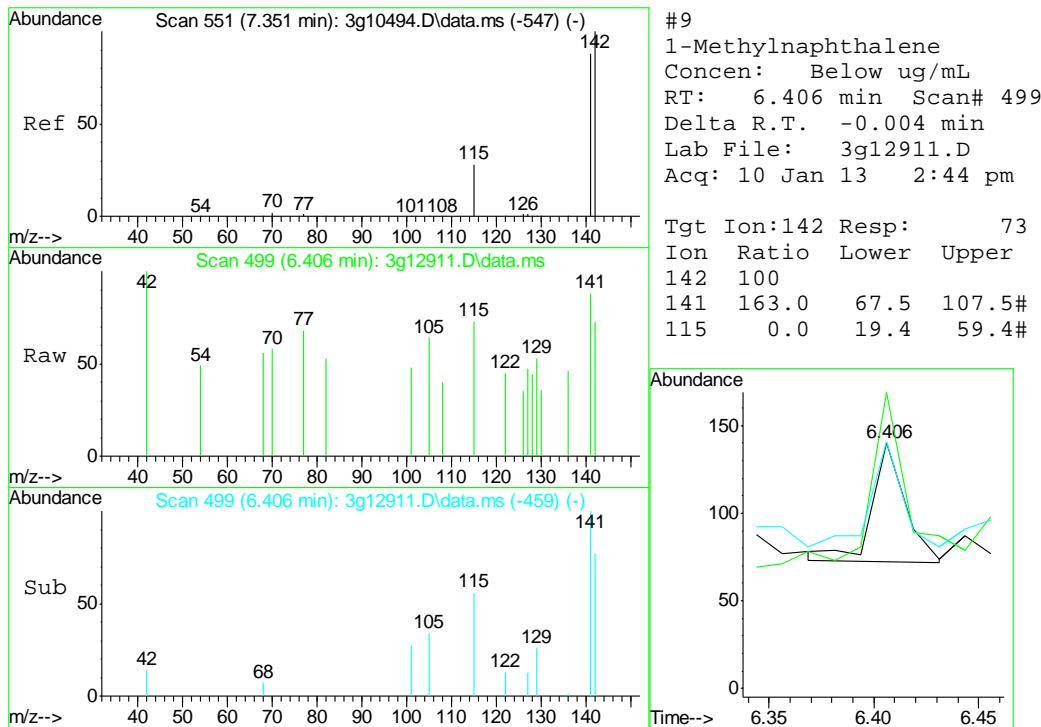


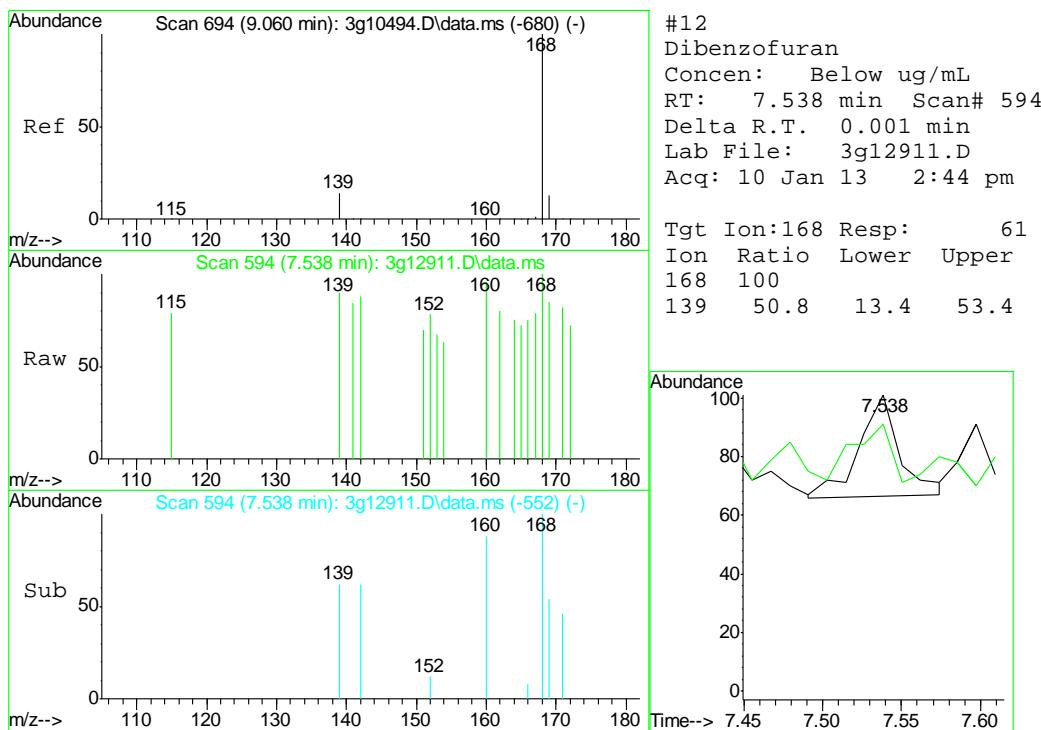
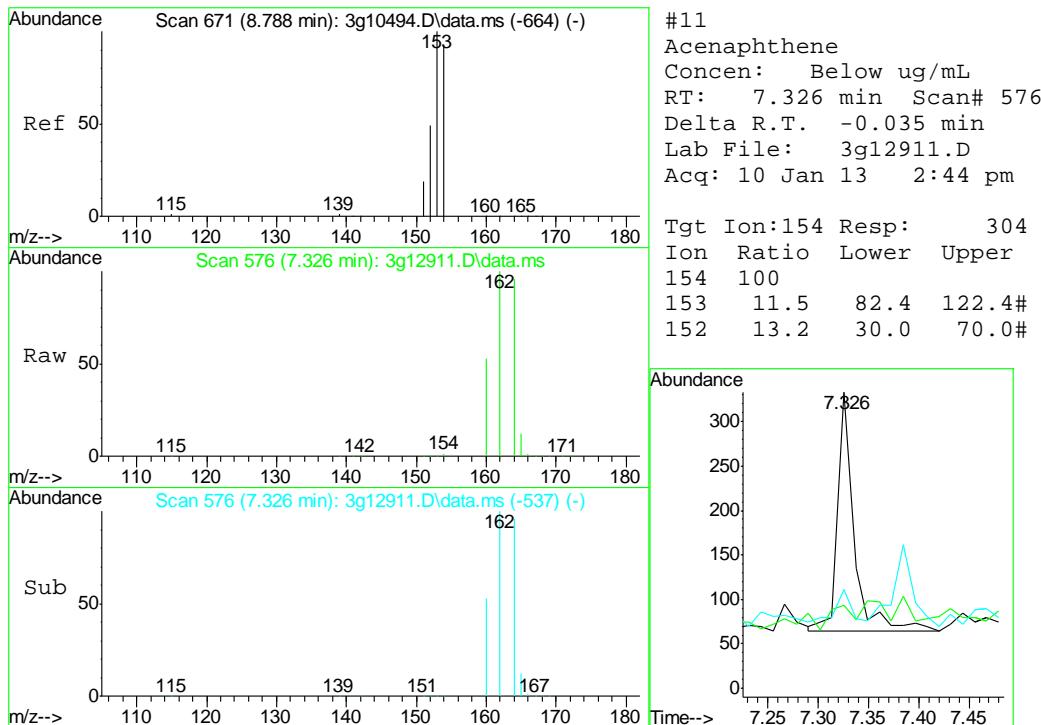


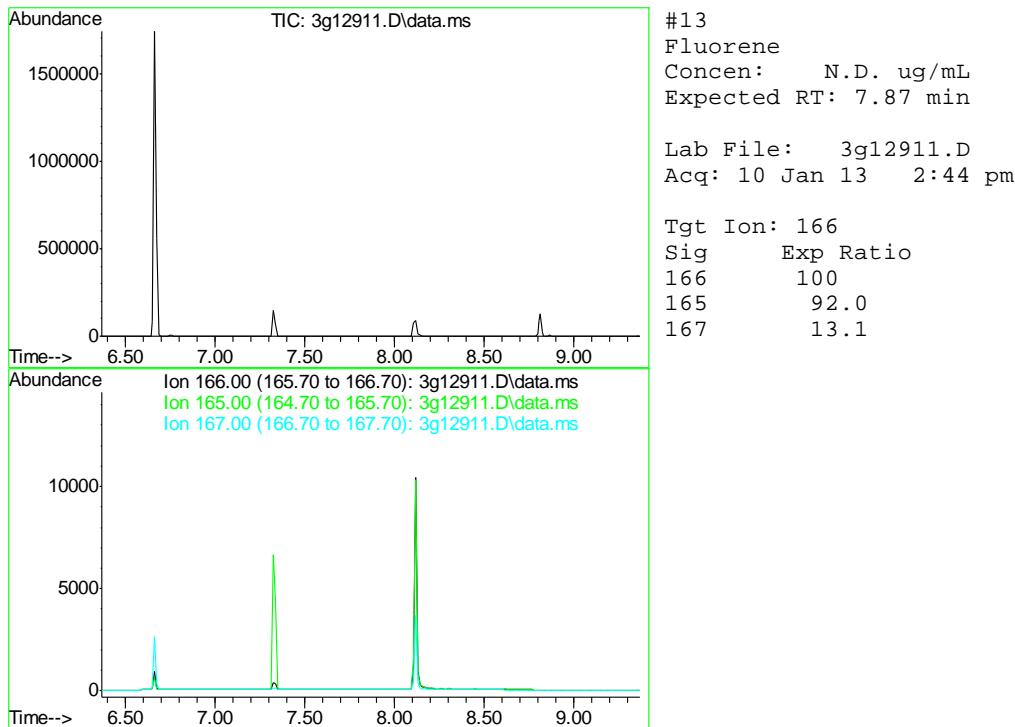






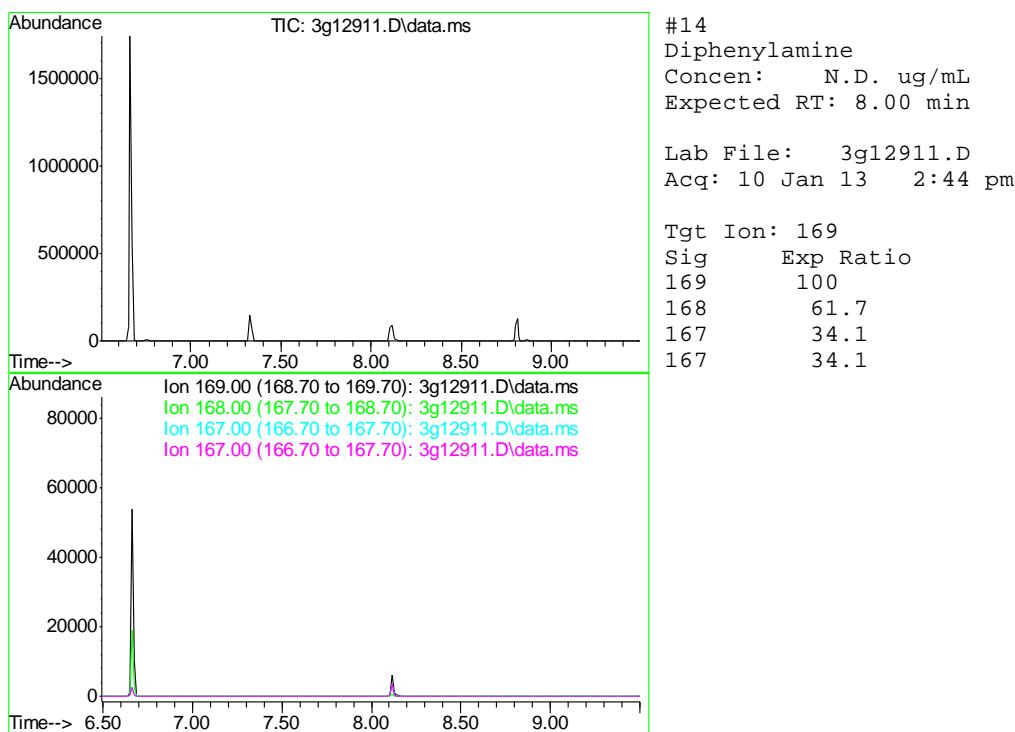


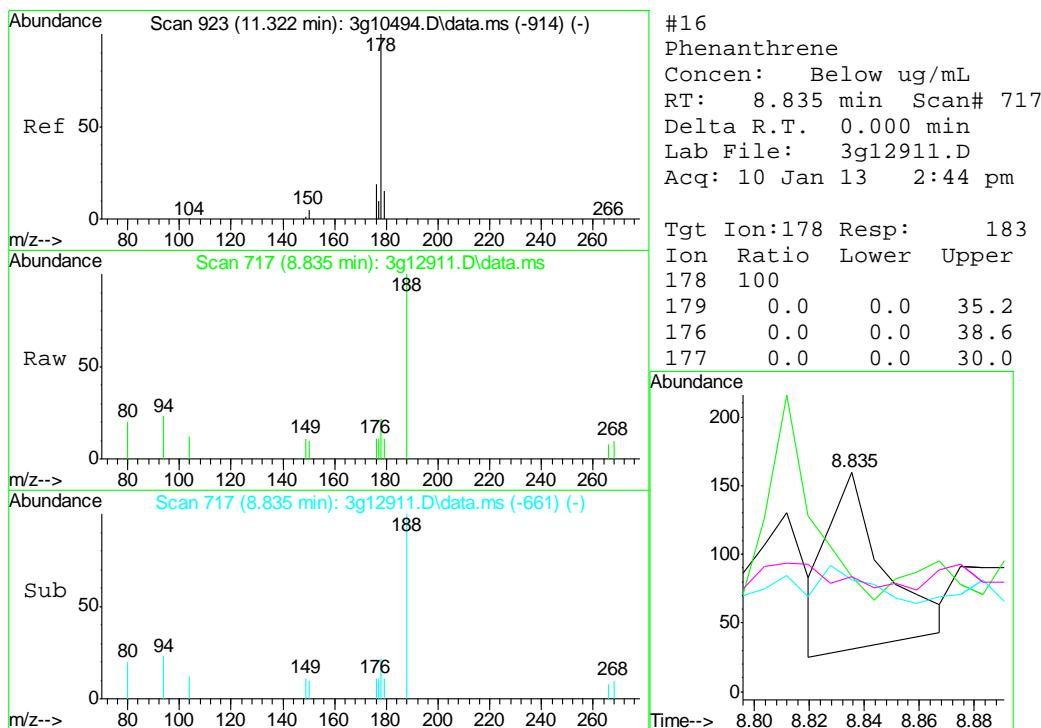
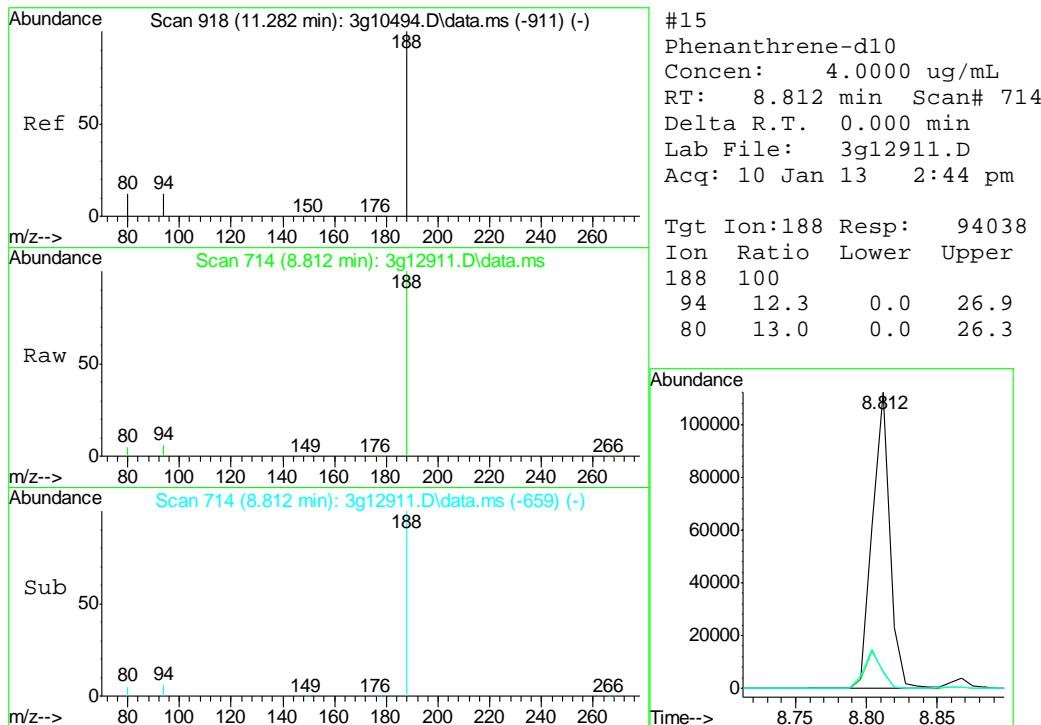


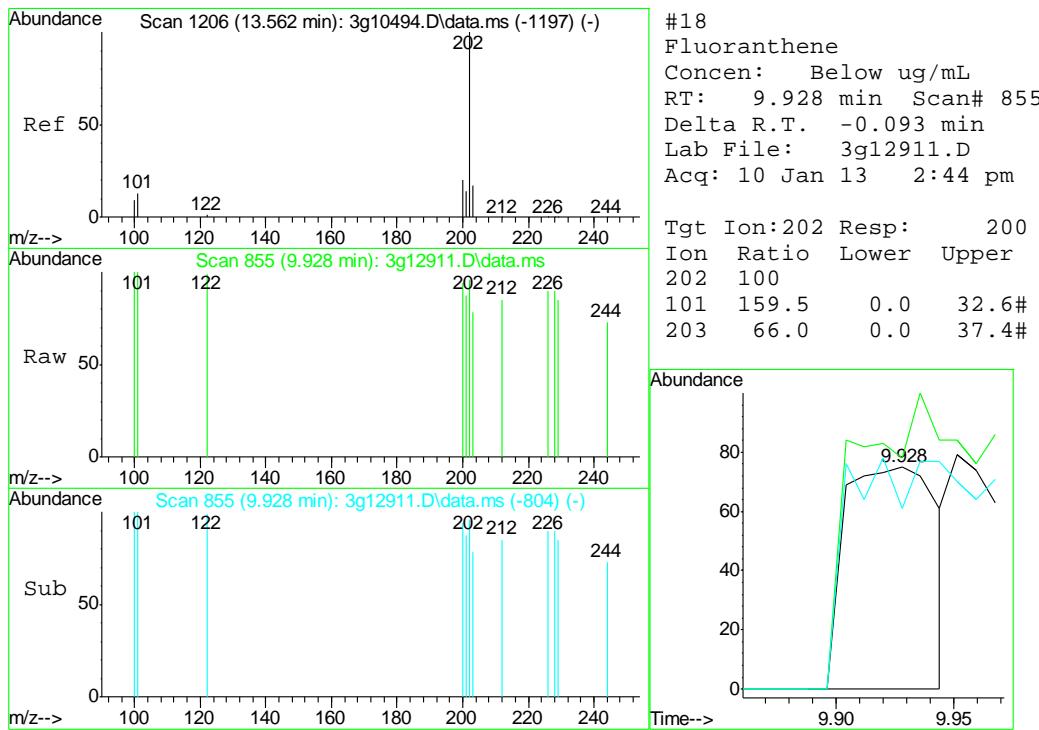
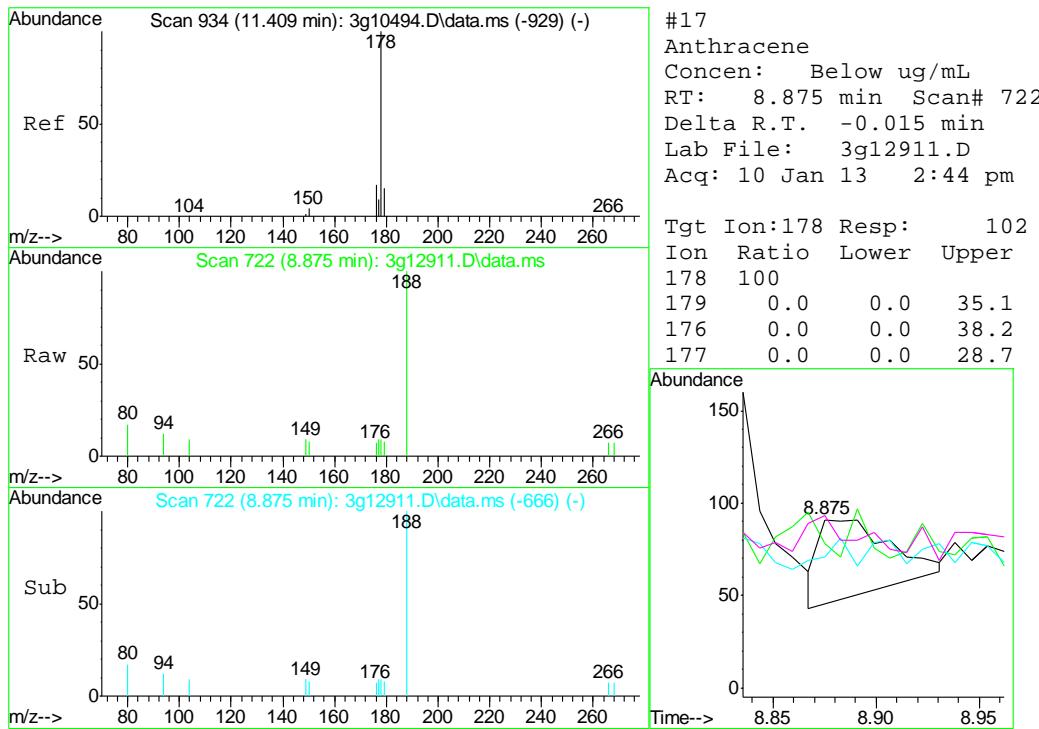


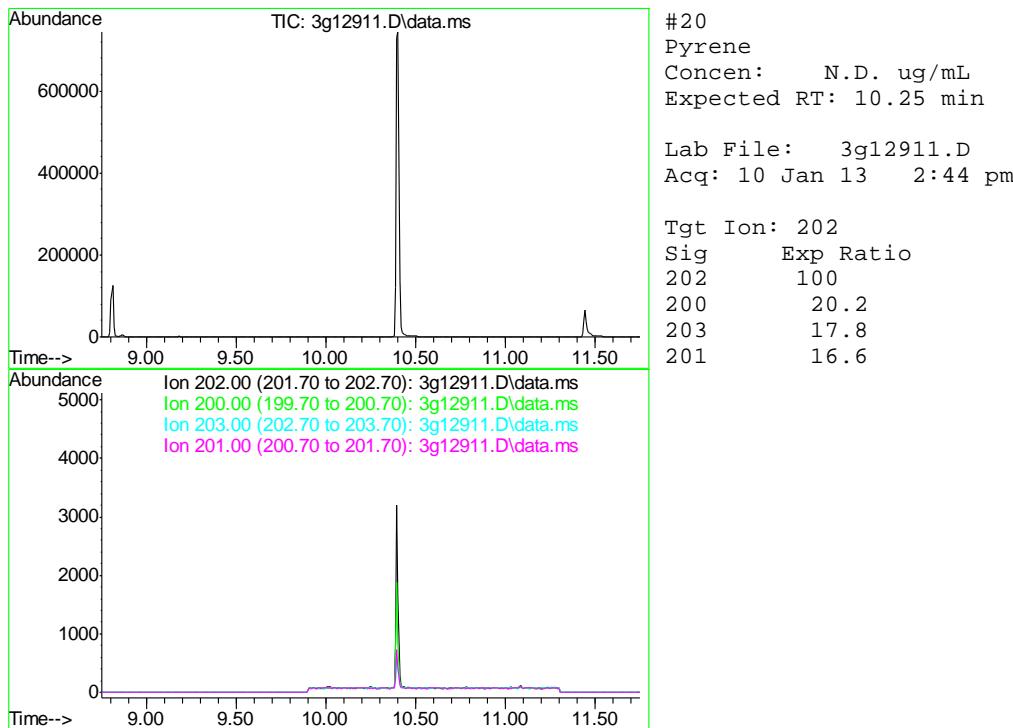
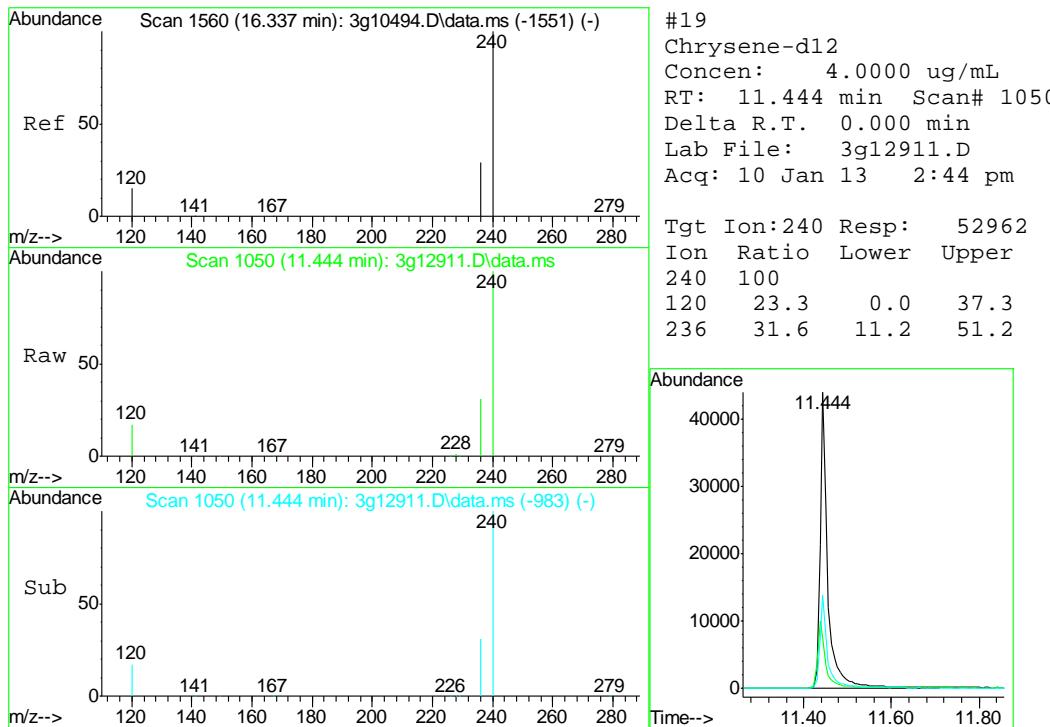
9.2.1

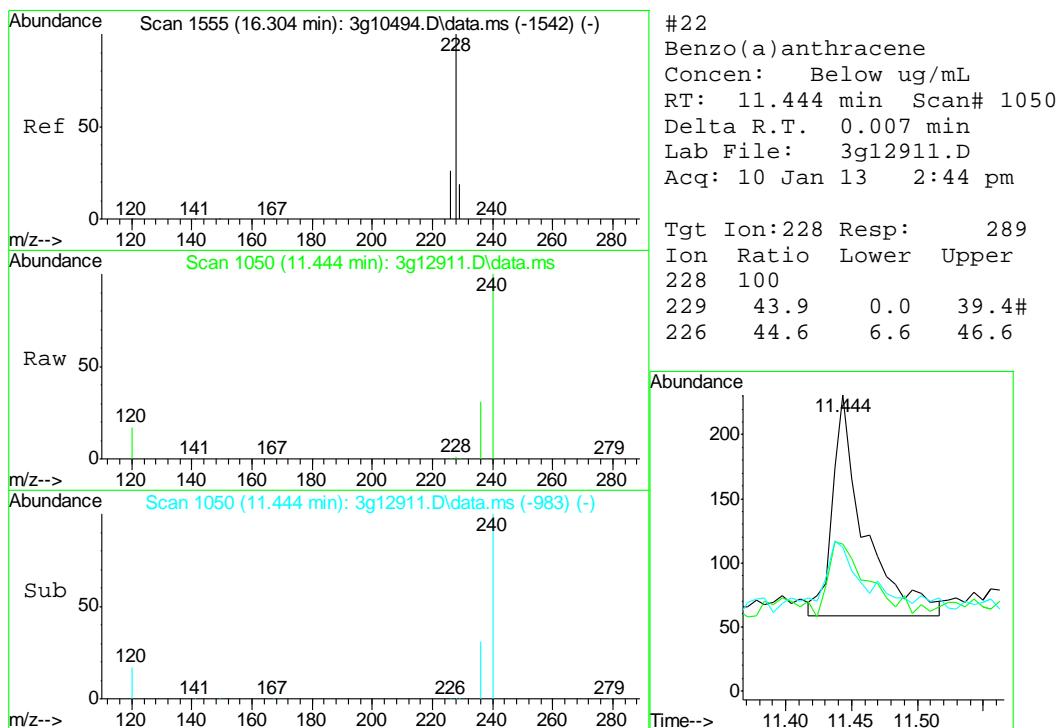
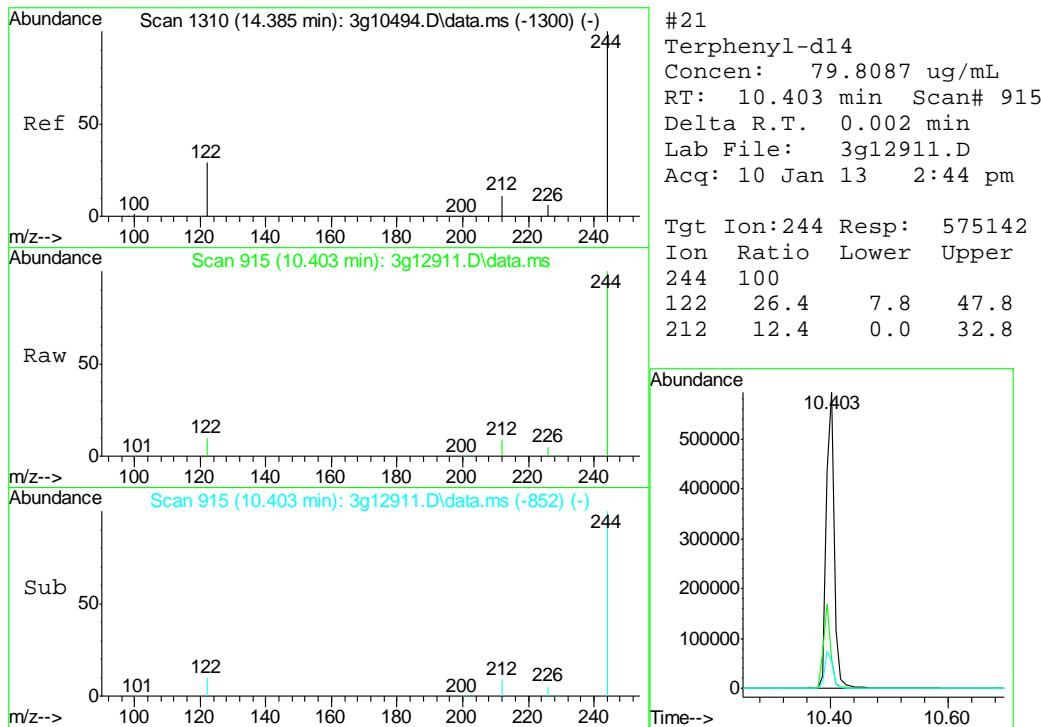
9

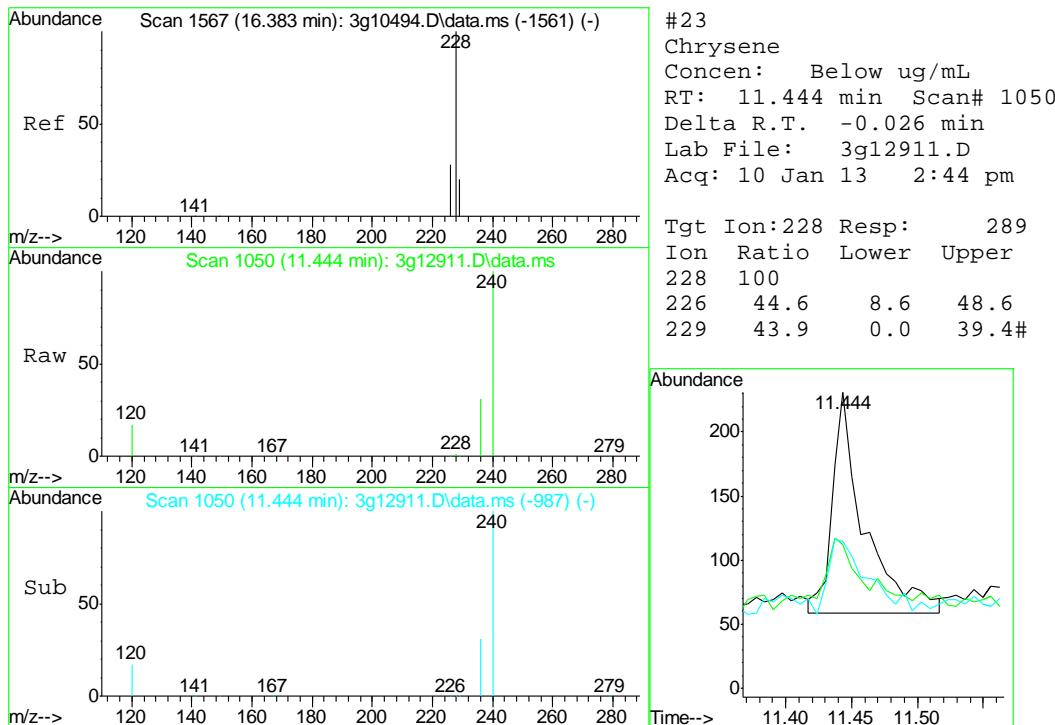






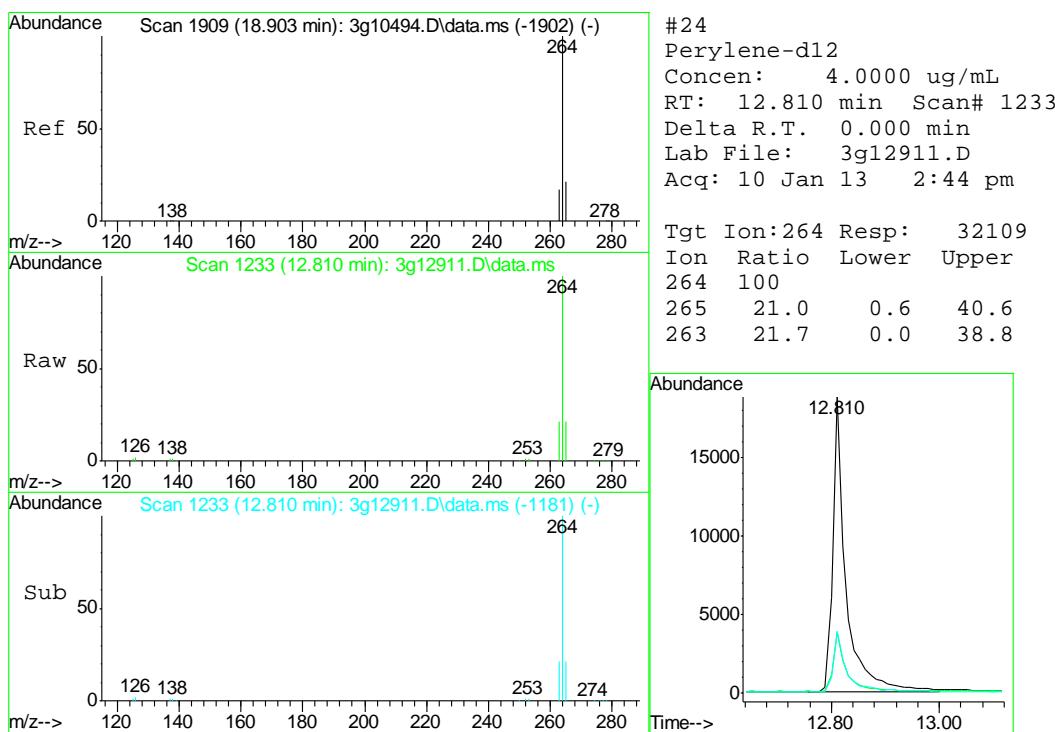


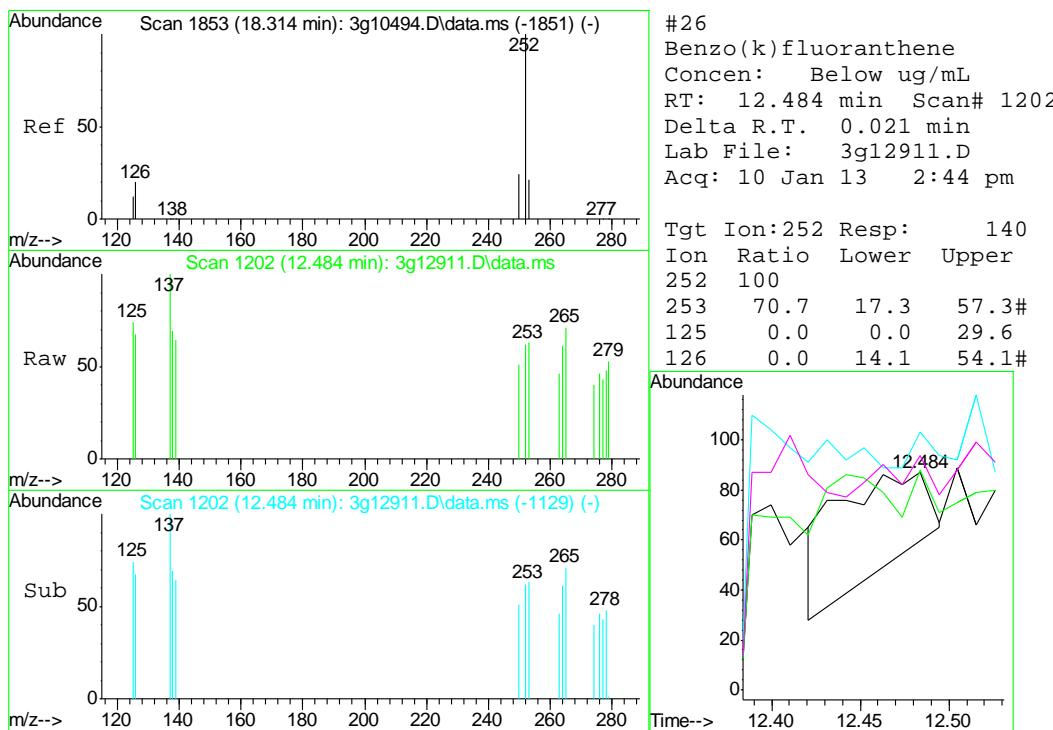
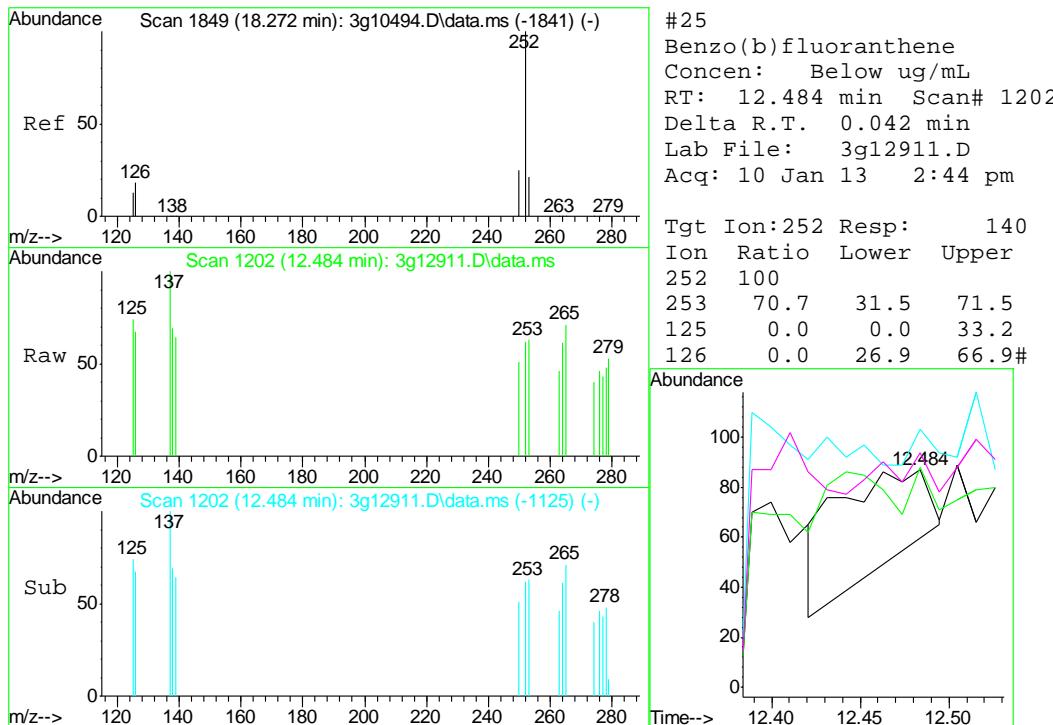


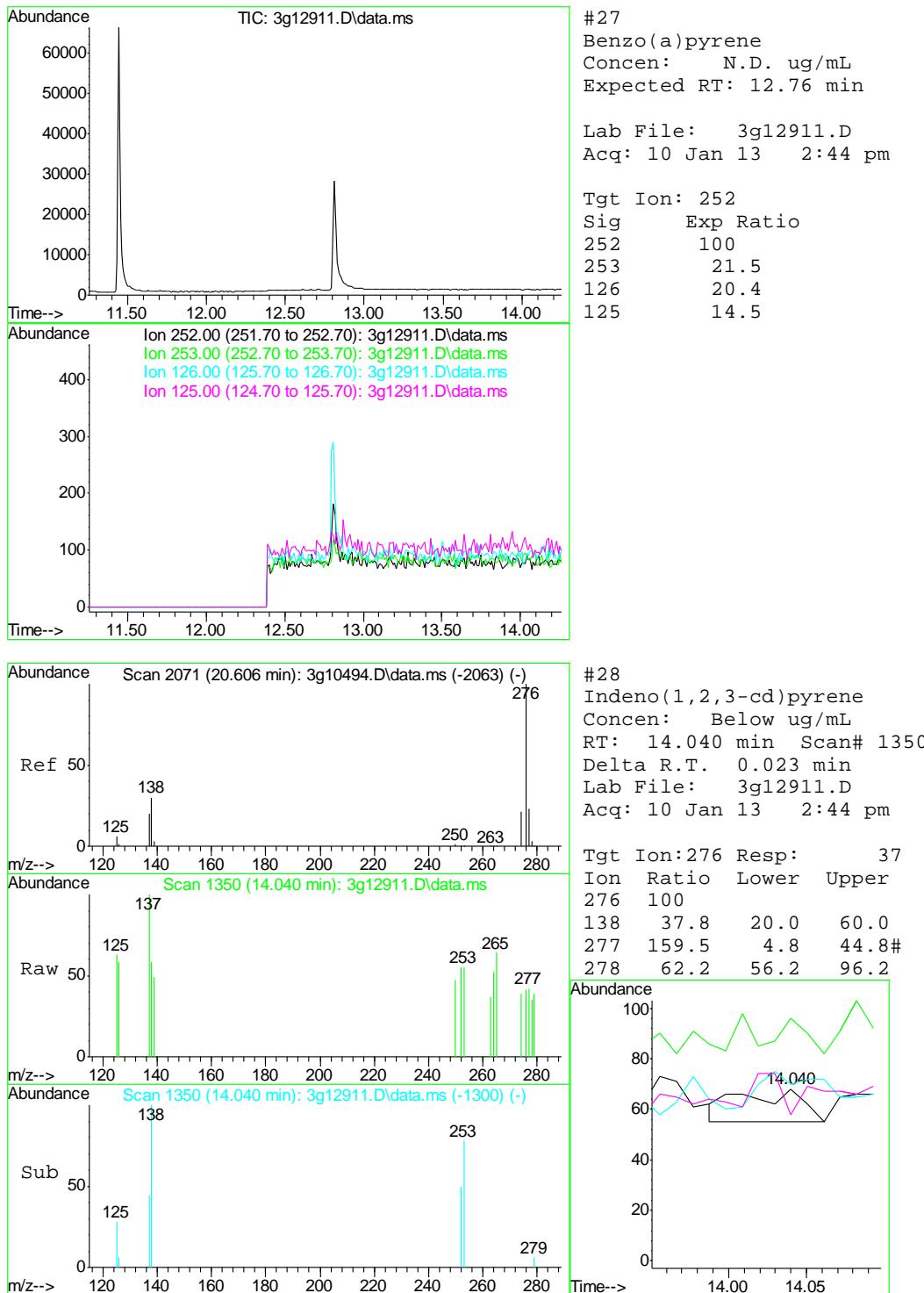


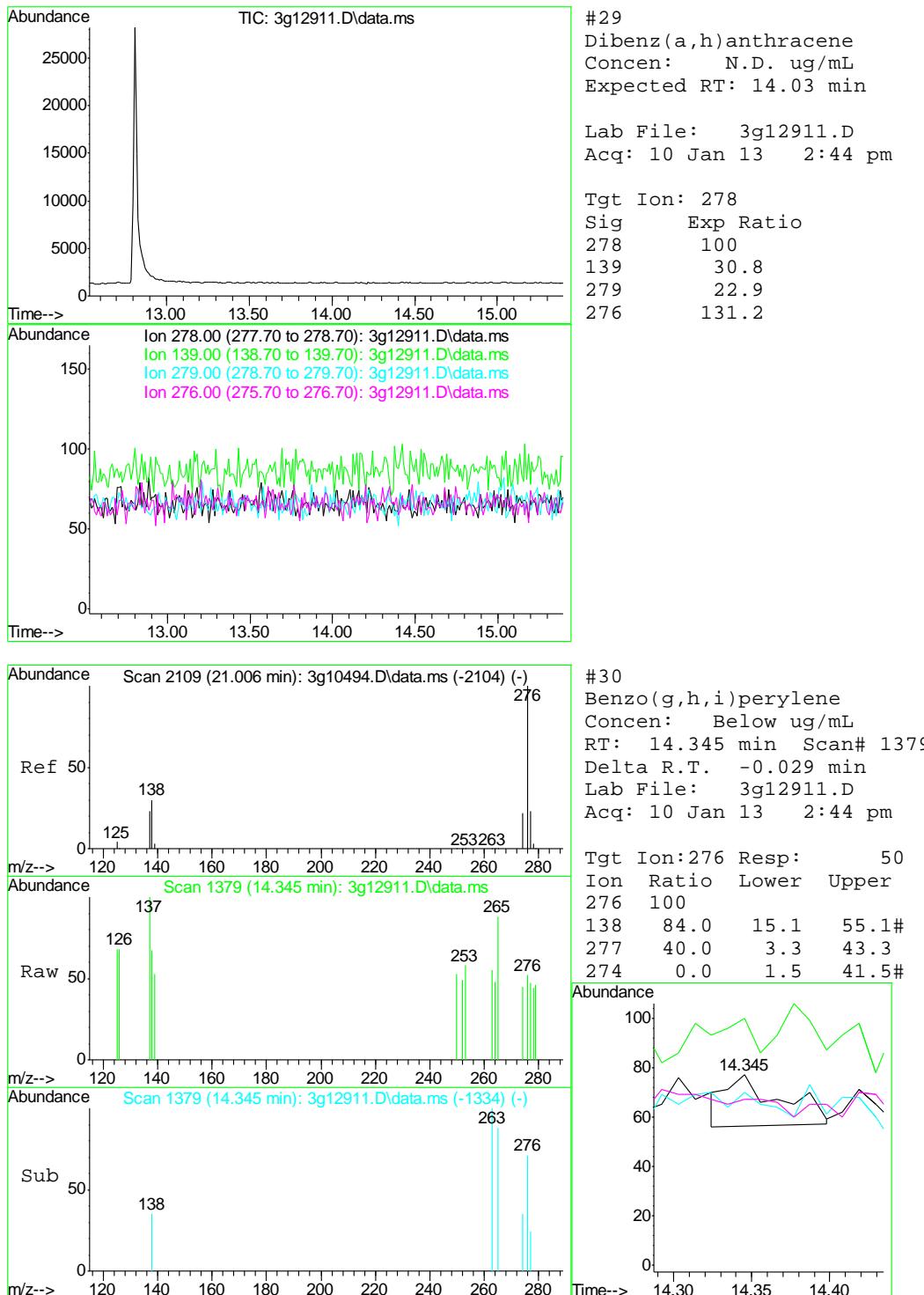
9.2.1

9











## GC Volatiles

---

### QC Data Summaries

---

Includes the following where applicable:

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

**Method Blank Summary**

**Job Number:** D42316  
**Account:** XTOKWR XTO Energy  
**Project:** XTO Love Ranch 8

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
GGB1042-MB	GB19076.D	1	01/07/13	SK	n/a	n/a	GGB1042

The QC reported here applies to the following samples:

**Method:** SW846 8015B

D42316-1

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	85% 60-140%

10.1.1

10

## Blank Spike Summary

Page 1 of 1

Job Number: D42316

Account: XTOKWR XTO Energy

Project: XTO Love Ranch 8

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
GGB1042-BS	GB19077.D	1	01/07/13	SK	n/a	n/a	GGB1042

The QC reported here applies to the following samples:

Method: SW846 8015B

D42316-1

CAS No.	Compound	Spike mg/kg	BSP mg/kg	BSP %	Limits
	TPH-GRO (C6-C10)	110	118	107	70-130

CAS No.	Surrogate Recoveries	BSP	Limits
120-82-1	1,2,4-Trichlorobenzene	102%	60-140%

10.2.1  
**10**

---

\* = Outside of Control Limits.

# Matrix Spike/Matrix Spike Duplicate Summary

Page 1 of 1

Job Number: D42316

Account: XTOKWR XTO Energy

Project: XTO Love Ranch 8

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
D42317-1MS	GB19079.D	1	01/07/13	SK	n/a	n/a	GGB1042
D42317-1MSD	GB19080.D	1	01/07/13	SK	n/a	n/a	GGB1042
D42317-1	GB19078.D	1	01/07/13	SK	n/a	n/a	GGB1042

The QC reported here applies to the following samples:

Method: SW846 8015B

D42316-1

CAS No.	Compound	D42317-1		Spike	MS	MS	MSD	MSD	RPD	Limits Rec/RPD
		mg/kg	Q	mg/kg	mg/kg	%	mg/kg	%		
	TPH-GRO (C6-C10)	ND		144	158	110	155	108	2	70-130/30

CAS No.	Surrogate Recoveries	MS	MSD	D42317-1	Limits
120-82-1	1,2,4-Trichlorobenzene	97%	94%	89%	60-140%

\* = Outside of Control Limits.

10.3.1  
10



## GC Volatiles

---

Raw Data

---

Judy Nelson  
 01/08/13 09:34

## Quantitation Report (QT Reviewed)

Signal #1 : Y:\1\DATA\010713\GB19081.D\FID1A.CH Vial: 8  
 Signal #2 : Y:\1\DATA\010713\GB19081.D\FID2B.CH  
 Acq On : 7 Jan 2013 4:03 pm Operator: StephK  
 Sample : D42316-1, 50X Inst : GC/MS Ins  
 Misc : GC3337,GGB1042,5.022,,100,5,1 Multiplr: 1.00  
 IntFile Signal #1: TVH1.E IntFile Signal #2: FB2.E  
 Quant Time: Jan 07 16:40:41 2013 Quant Results File: TB868GB868SOIL.RES

Quant Method : C:\MSDCHEM\1...\TB868GB868SOIL.M (Chemstation Integrator)  
 Title : 8015B/8021B TVH/BTEX  
 Last Update : Mon Jan 07 12:45:52 2013  
 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	14.36	3841859	122.610 %	m
10) S	1,2,4-Trichlorobenzene (P)	14.36	16537697	101.753 %	m

**Target Compounds**

1) H	TVH-Gasoline	7.23	133314210	2.088 mg/L
4) T	Methyl-t-butyl-ether	0.00	0	N.D. ug/L d
5) T	Benzene	4.16	83088	0.206 ug/L
6) T	Toluene	7.66	1266424	3.196 ug/L m
7) T	Ethylbenzene	10.28	1227942	3.630 ug/L
8) T	m,p-Xylene	10.46	25045415	68.245 ug/L
9) T	o-Xylene	10.96	4828944	14.706 ug/L
11) T	Naphthalene	14.54	7972179	40.405 ug/L m

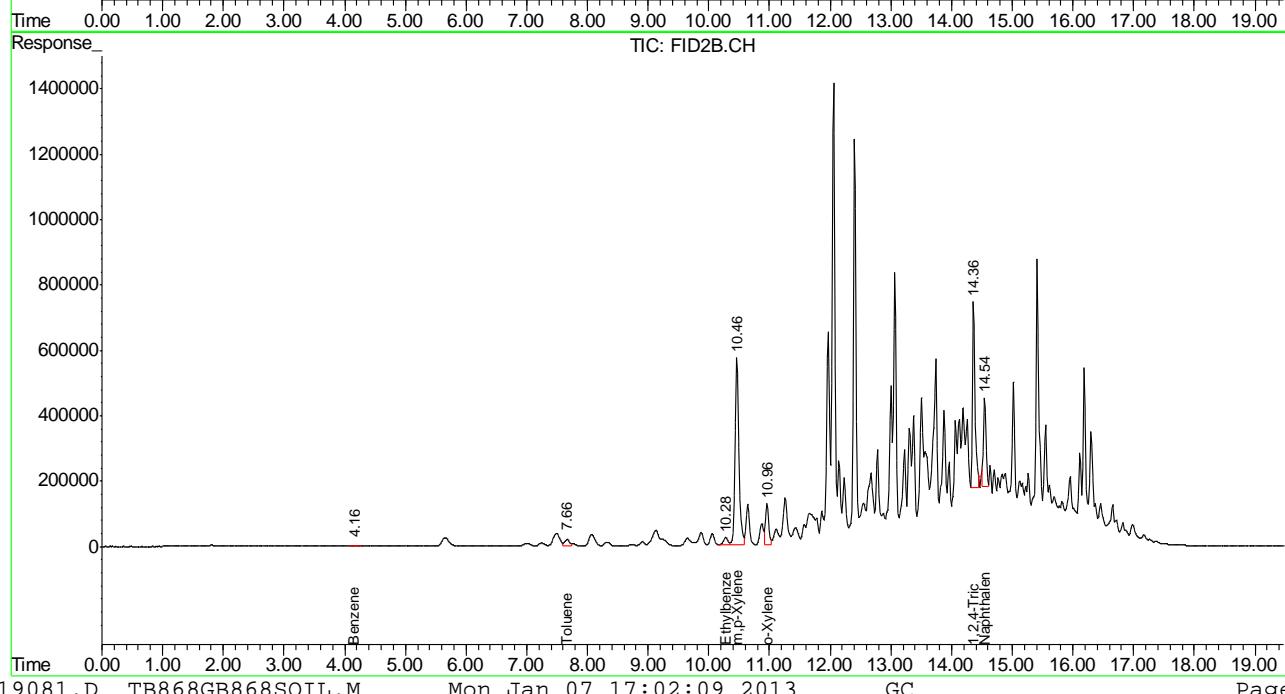
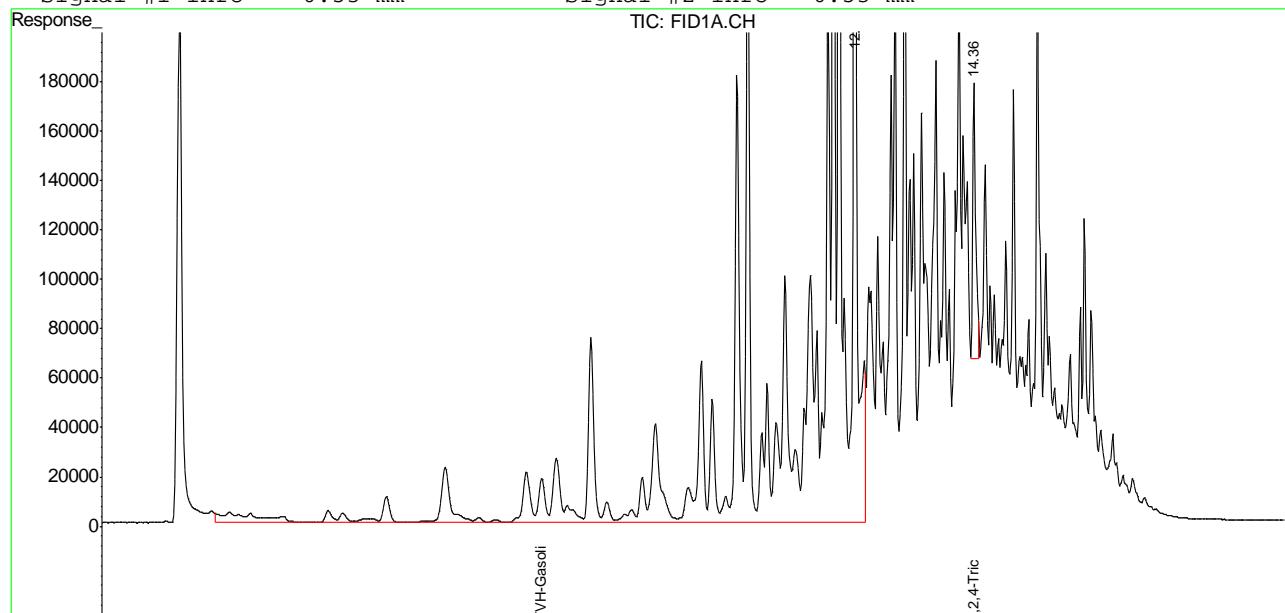
(f)=RT Delta > 1/2 Window (m)=manual int.  
 GB19081.D TB868GB868SOIL.M Mon Jan 07 17:02:09 2013 GC

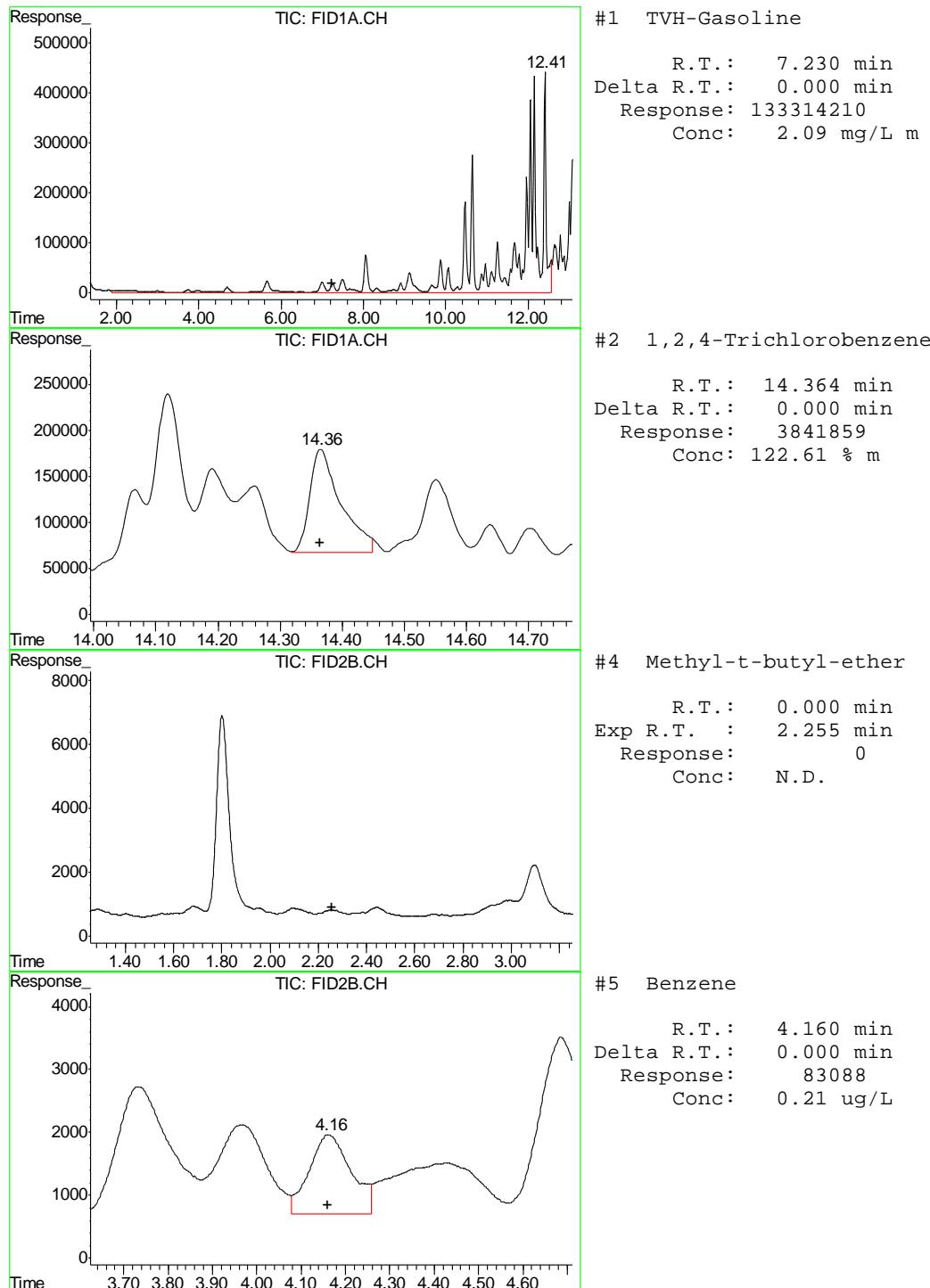
## Quantitation Report (QT Reviewed)

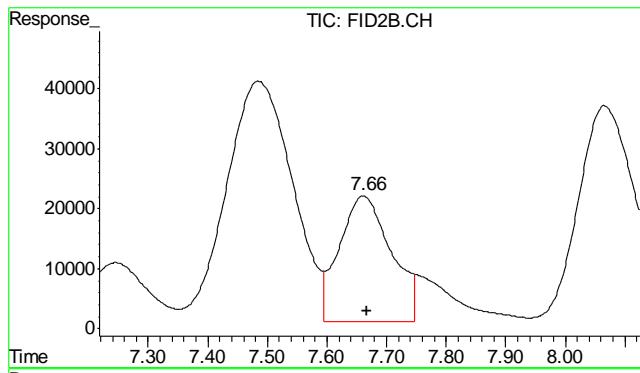
Signal #1 : Y:\1\DATA\010713\GB19081.D\FID1A.CH Vial: 8  
 Signal #2 : Y:\1\DATA\010713\GB19081.D\FID2B.CH  
 Acq On : 7 Jan 2013 4:03 pm Operator: StephK  
 Sample : D42316-1, 50X Inst : GC/MS Ins  
 Misc : GC3337,GGB1042,5.022,,100,5,1 Multiplr: 1.00  
 IntFile Signal #1: TVH1.E IntFile Signal #2: FB2.E  
 Quant Time: Jan 7 16:40 2013 Quant Results File: TB868GB868SOIL.RES

Quant Method : C:\MSDCHEM\1...\TB868GB868SOIL.M (Chemstation Integrator)  
 Title : 8015B/8021B TVH/BTEX  
 Last Update : Mon Jan 07 12:45:52 2013  
 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

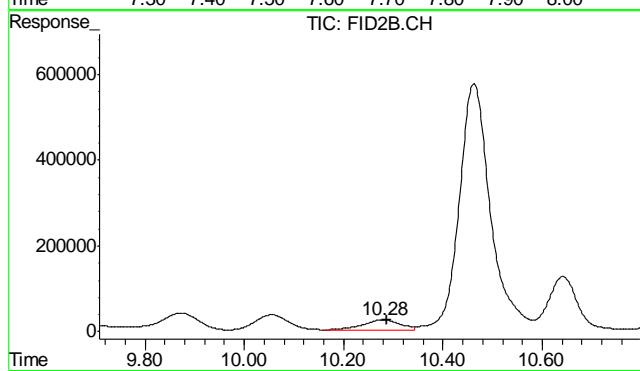






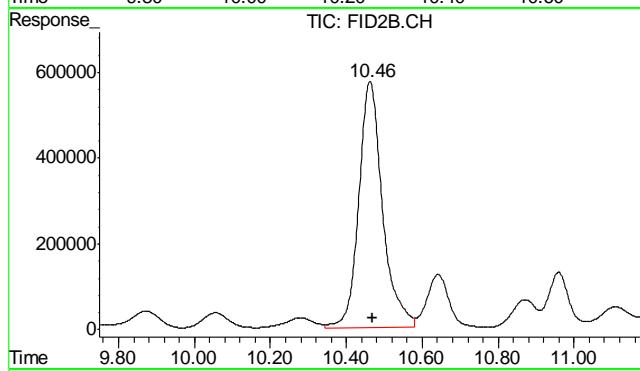
#6 Toluene

R.T.: 7.661 min  
Delta R.T.: -0.005 min  
Response: 1266424  
Conc: 3.20 ug/L m



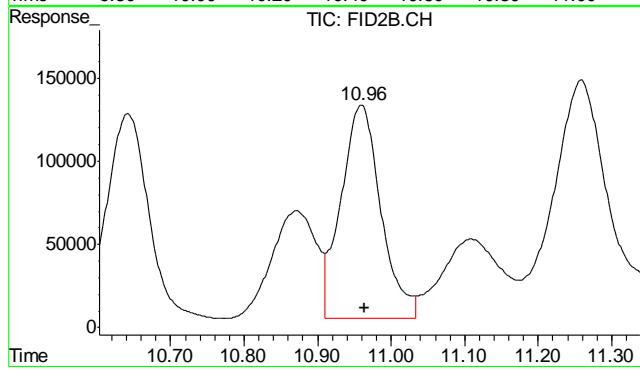
#7 Ethylbenzene

R.T.: 10.279 min  
Delta R.T.: -0.009 min  
Response: 1227942  
Conc: 3.63 ug/L



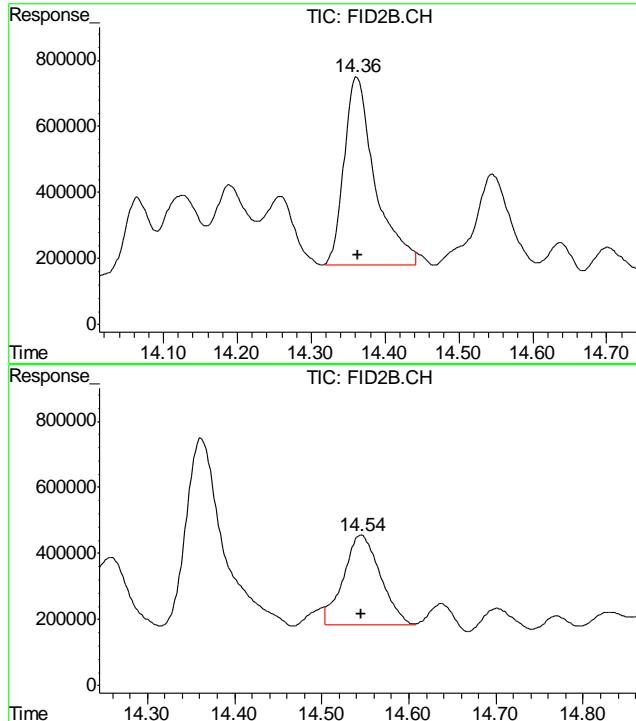
#8 m,p-Xylene

R.T.: 10.463 min  
Delta R.T.: -0.005 min  
Response: 25045415  
Conc: 68.25 ug/L

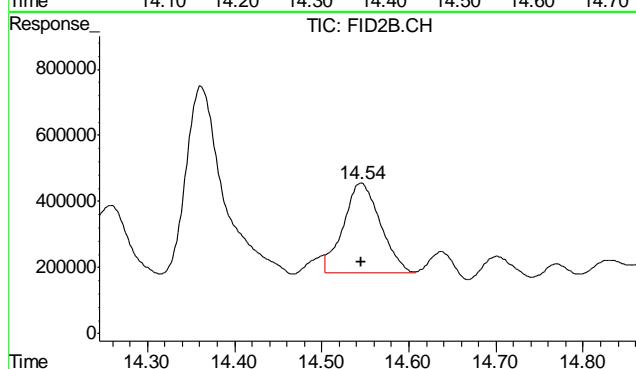


#9 o-Xylene

R.T.: 10.960 min  
Delta R.T.: -0.004 min  
Response: 4828944  
Conc: 14.71 ug/L



#10 1,2,4-Trichlorobenzene (P)  
 R.T.: 14.361 min  
 Delta R.T.: -0.002 min  
 Response: 16537697  
 Conc: 101.75 % m



#11 Naphthalene  
 R.T.: 14.545 min  
 Delta R.T.: 0.000 min  
 Response: 7972179  
 Conc: 40.40 ug/L m

11.1.1

**Manual Integrations  
APPROVED  
(compounds with "m" flag)**  
**Judy Nelson  
01/08/13 09:34**

## Quantitation Report (QT Reviewed)

Signal #1 : Y:\1\DATA\010713\GB19076.D\FID1A.CH Vial: 3  
 Signal #2 : Y:\1\DATA\010713\GB19076.D\FID2B.CH  
 Acq On : 7 Jan 2013 1:06 pm Operator: StephK  
 Sample : MB Inst : GC/MS Ins  
 Misc : GC3337,GGB1042,5.000,,100,5,1 Multiplr: 1.00  
 IntFile Signal #1: TVH1.E IntFile Signal #2: FB2.E  
 Quant Time: Jan 07 14:37:39 2013 Quant Results File: TB868GB868SOIL.RES

Quant Method : C:\MSDCHEM\1...\TB868GB868SOIL.M (Chemstation Integrator)  
 Title : 8015B/8021B TVH/BTEX  
 Last Update : Mon Jan 07 12:45:52 2013  
 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	14.37	2667492	85.131 %
10) S	1,2,4-Trichlorobenzene (P)	14.37	13776254	84.763 % m

**Target Compounds**

1) H	TVH-Gasoline	7.23	3000971	<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	7.68	134125	0.338 ug/L
7) T	Ethylbenzene	0.00	0	N.D. ug/L d
8) T	m,p-Xylene	0.00	0	N.D. ug/L d
9) T	o-Xylene	0.00	0	N.D. ug/L d
11) T	Naphthalene	14.55	24733	0.125 ug/L m

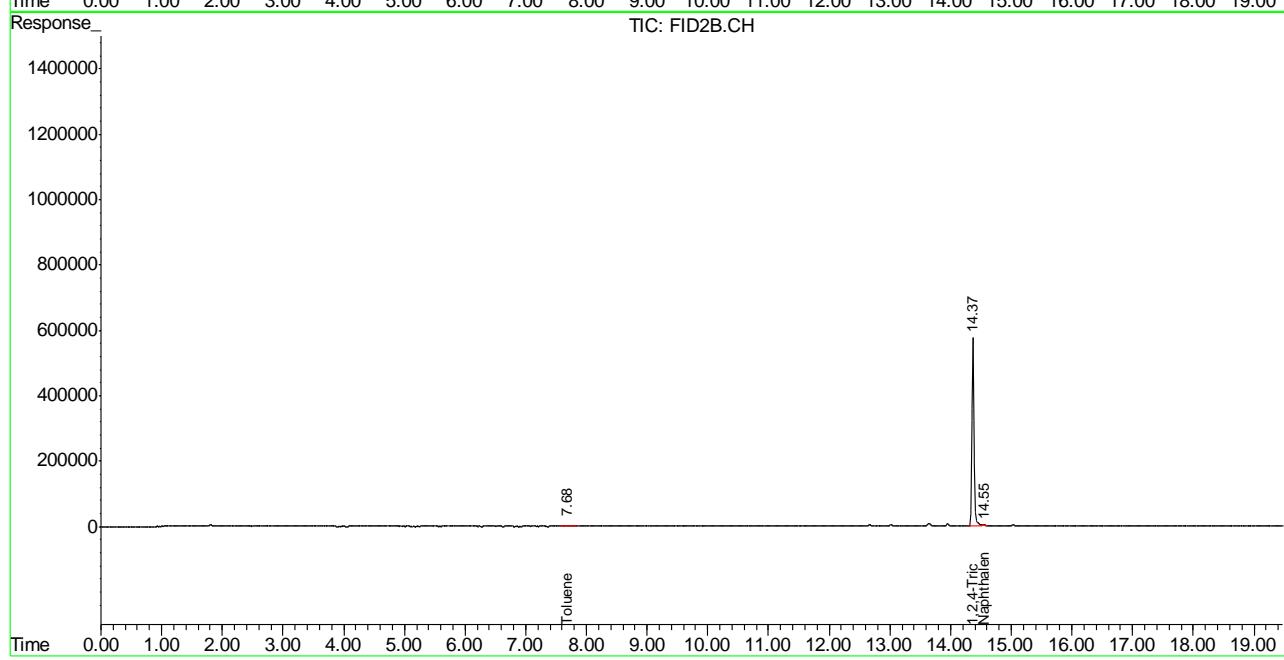
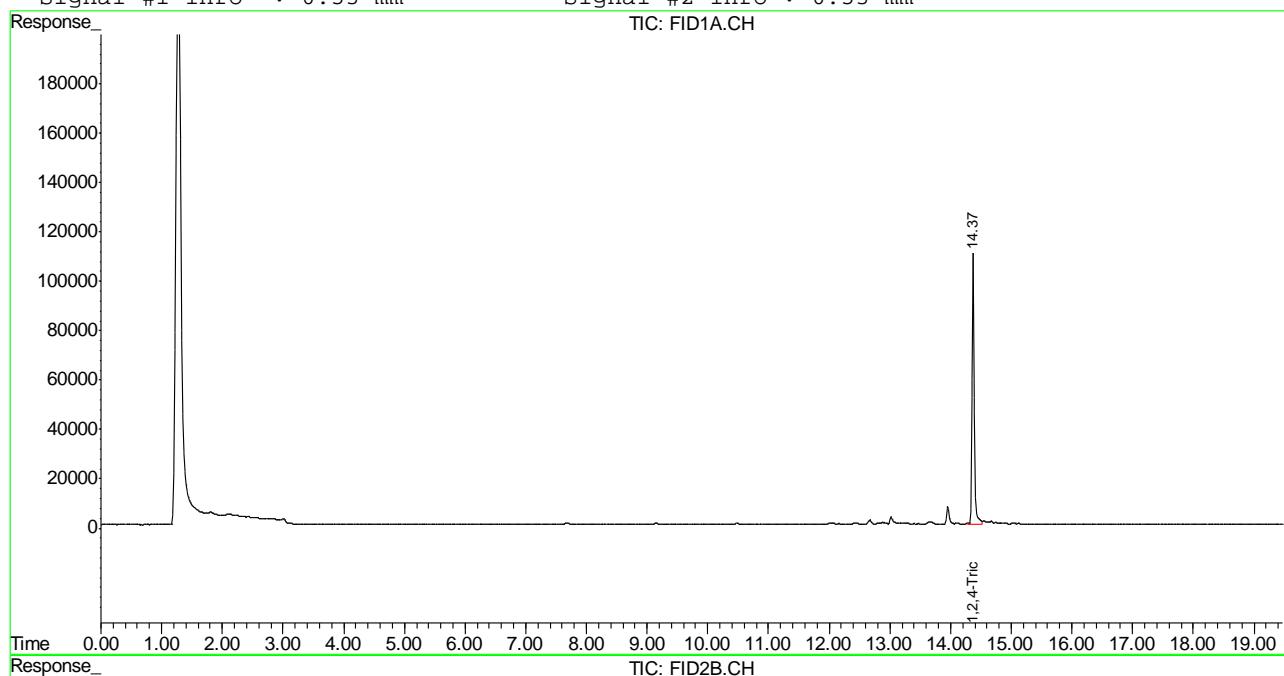
(f)=RT Delta > 1/2 Window (m)=manual int.  
 GB19076.D TB868GB868SOIL.M Mon Jan 07 17:01:54 2013 GC

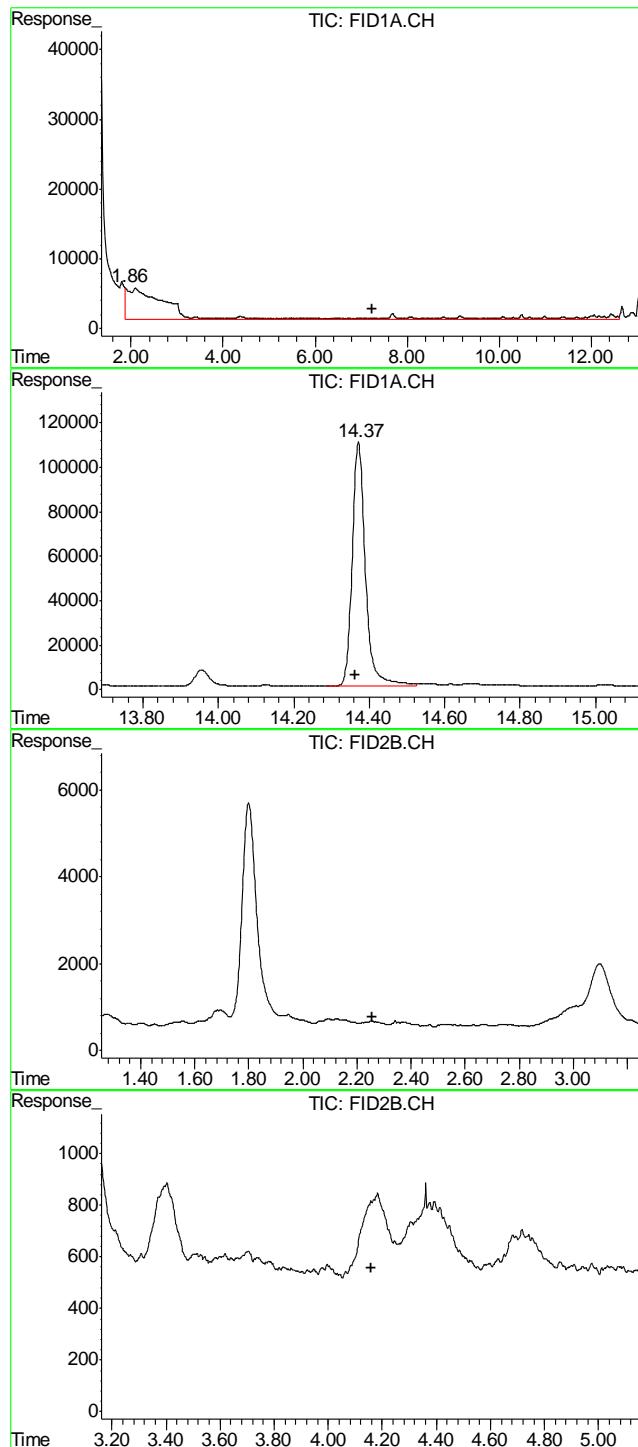
## Quantitation Report (QT Reviewed)

Signal #1 : Y:\1\DATA\010713\GB19076.D\FID1A.CH Vial: 3  
 Signal #2 : Y:\1\DATA\010713\GB19076.D\FID2B.CH  
 Acq On : 7 Jan 2013 1:06 pm Operator: StephK  
 Sample : MB Inst : GC/MS Ins  
 Misc : GC3337,GGB1042,5.000,,100,5,1 Multiplr: 1.00  
 IntFile Signal #1: TVH1.E IntFile Signal #2: FB2.E  
 Quant Time: Jan 7 14:39 2013 Quant Results File: TB868GB868SOIL.RES

Quant Method : C:\MSDCHEM\1...\TB868GB868SOIL.M (Chemstation Integrator)  
 Title : 8015B/8021B TVH/BTEX  
 Last Update : Mon Jan 07 12:45:52 2013  
 Response via : Multiple Level Calibration  
 DataAcq Meth : TVB4.M

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



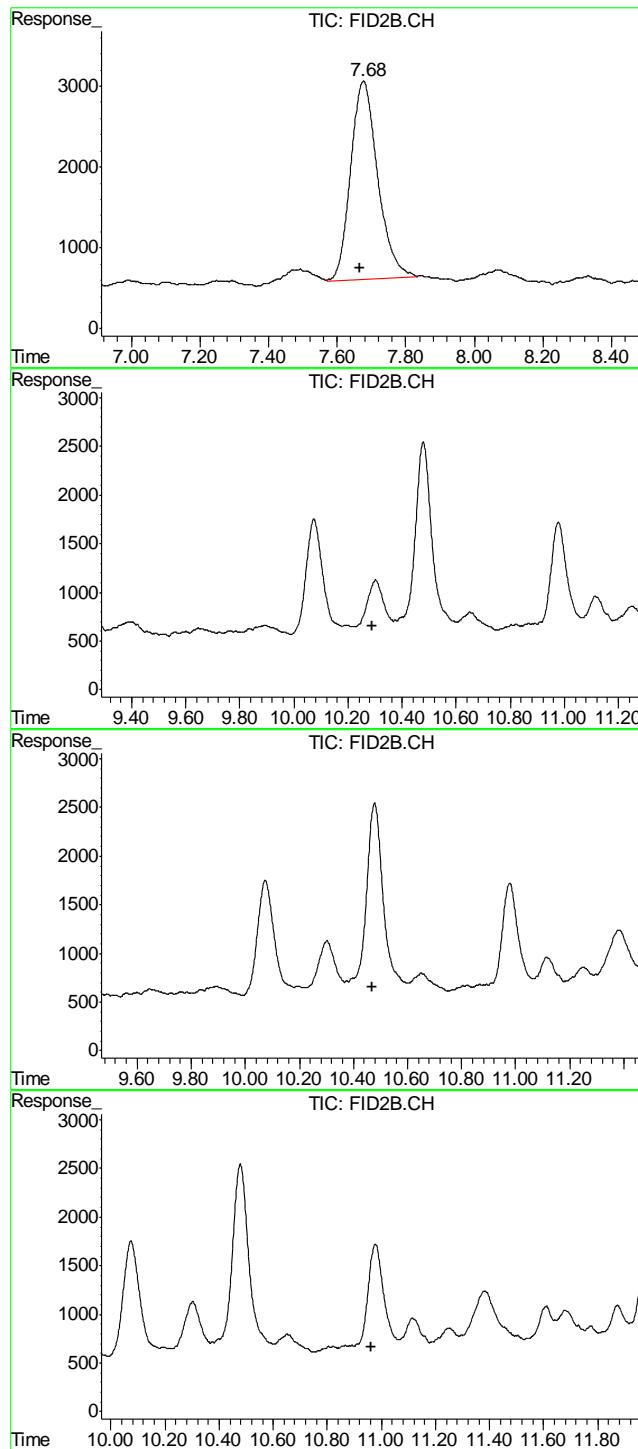


#1 TVH-Gasoline  
 R.T.: 7.230 min  
 Delta R.T.: 0.000 min  
 Response: 3000971  
 Conc: N.D.

#2 1,2,4-Trichlorobenzene  
 R.T.: 14.372 min  
 Delta R.T.: 0.007 min  
 Response: 2667492  
 Conc: 85.13 %

#4 Methyl-t-butyl-ether  
 R.T.: 0.000 min  
 Exp R.T. : 2.255 min  
 Response: 0  
 Conc: N.D.

#5 Benzene  
 R.T.: 0.000 min  
 Exp R.T. : 4.161 min  
 Response: 0  
 Conc: N.D.

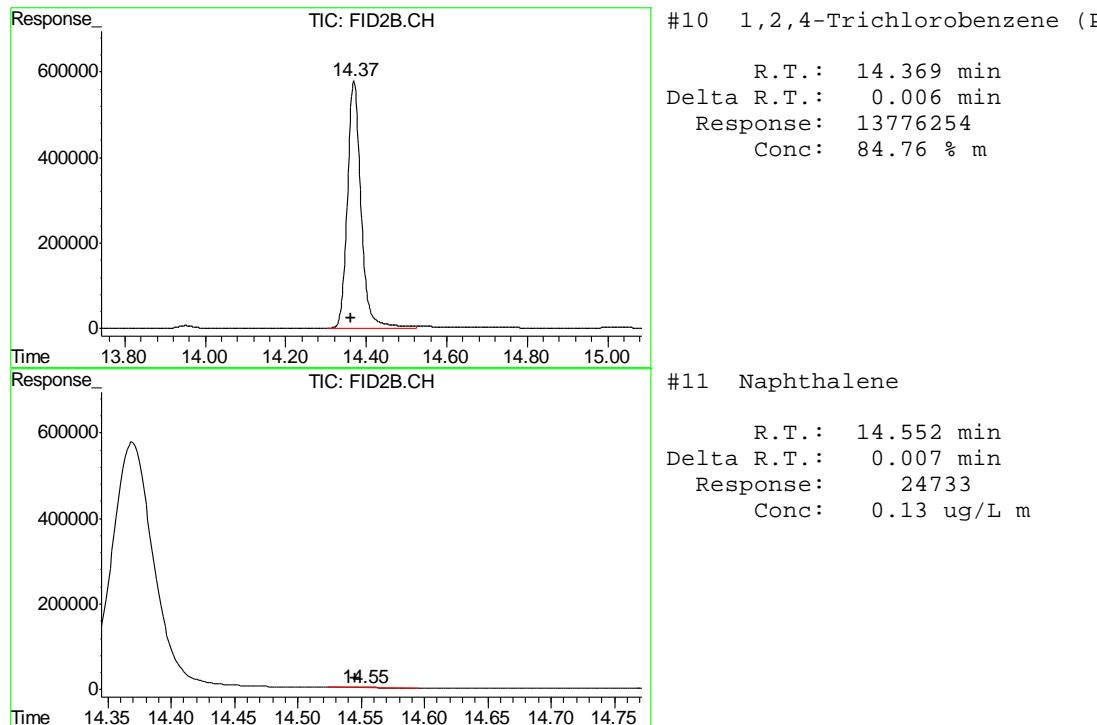


#6 Toluene  
R.T.: 7.677 min  
Delta R.T.: 0.011 min  
Response: 134125  
Conc: 0.34 ug/L

#7 Ethylbenzene  
R.T.: 0.000 min  
Exp R.T.: 10.288 min  
Response: 0  
Conc: N.D.

#8 m,p-Xylene  
R.T.: 0.000 min  
Exp R.T.: 10.468 min  
Response: 0  
Conc: N.D.

#9 o-Xylene  
R.T.: 0.000 min  
Exp R.T.: 10.963 min  
Response: 0  
Conc: N.D.





## GC Semi-volatiles

---

### QC Data Summaries

---

Includes the following where applicable:

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

**Method Blank Summary**

Page 1 of 1

Job Number: D42316

Account: XTOKWR XTO Energy

Project: XTO Love Ranch 8

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
OP7201-MB	FD21006.D	1	01/09/13	AV	01/09/13	OP7201	GFD1058

The QC reported here applies to the following samples:

**Method:** SW846-8015B

D42316-1

CAS No.	Compound	Result	RL	MDL	Units	Q
	TPH-DRO (C10-C28)	ND	6.7	4.0	mg/kg	

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

## Blank Spike Summary

Page 1 of 1

Job Number: D42316

Account: XTOKWR XTO Energy

Project: XTO Love Ranch 8

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
OP7201-BS	FD21008.D	1	01/09/13	AV	01/09/13	OP7201	GFD1058

The QC reported here applies to the following samples:

Method: SW846-8015B

D42316-1

CAS No.	Compound	Spike mg/kg	BSP mg/kg	BSP %	Limits
	TPH-DRO (C10-C28)	667	597	90	48-130

CAS No.	Surrogate Recoveries	BSP	Limits
84-15-1	o-Terphenyl	83%	35-130%

\* = Outside of Control Limits.

12.2.1

12

# Matrix Spike/Matrix Spike Duplicate Summary

Page 1 of 1

Job Number: D42316

Account: XTOKWR XTO Energy

Project: XTO Love Ranch 8

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
OP7201-MS	FD21068.D	10	01/10/13	AV	01/09/13	OP7201	GFD1060
OP7201-MSD	FD21070.D	10	01/10/13	AV	01/09/13	OP7201	GFD1060
D42316-1	FD21072.D	10	01/10/13	AV	01/09/13	OP7201	GFD1060

The QC reported here applies to the following samples:

Method: SW846-8015B

D42316-1

CAS No.	Compound	D42316-1		Spike	MS	MS	MSD	MSD	RPD	Limits Rec/RPD
		mg/kg	Q	mg/kg	mg/kg	%	mg/kg	%		
	TPH-DRO (C10-C28)	5850		1050	5460	-37* a	6050	19* a	10	20-168/30

CAS No.	Surrogate Recoveries	MS	MSD	D42316-1	Limits
84-15-1	o-Terphenyl	58%	72%	69%	35-130%

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

\* = Outside of Control Limits.

12.3.1  
12



## GC Semi-volatiles

---

Raw Data

---

Manual Integrations  
APPROVED  
(compounds with "m" flag)

Judy Nelson
01/11/13 09:35

## Quantitation Report (QT Reviewed)

Data File : C:\MSDCHEM\2\DATA\2013\JAN\FD011013.SEC\FD21072.D Vial: 72  
 Acq On : 1-10-2013 09:05:02 PM Operator: ashleyv  
 Sample : D42316-1, 10x Inst : FID5  
 Misc : OP7201,GFD1060,30.06,,,1,10 Multiplr: 1.00  
 IntFile : autoint1.e  
 Quant Time: Jan 11 08:08:21 2013 Quant Results File: DRO-GFD983R.RES

Quant Method : C:\MSDCHEM\2...\DRO-GFD983R.M (Chemstation Integrator)  
 Title : 8015B TEH  
 Last Update : Tue Jan 08 09:15:06 2013  
 Response via : Initial Calibration  
 DataAcq Meth : DRODUAL.M

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

Compound	R.T.	Response	Conc Units
<hr/>			
System Monitoring Compounds			
1) S O-Terphenyl	8.96	7064675	138.172 mg/L m
<hr/>			
Target Compounds			
2) H TPH-DRO (c10-c28)	6.93	411428431	11132.163 mg/L

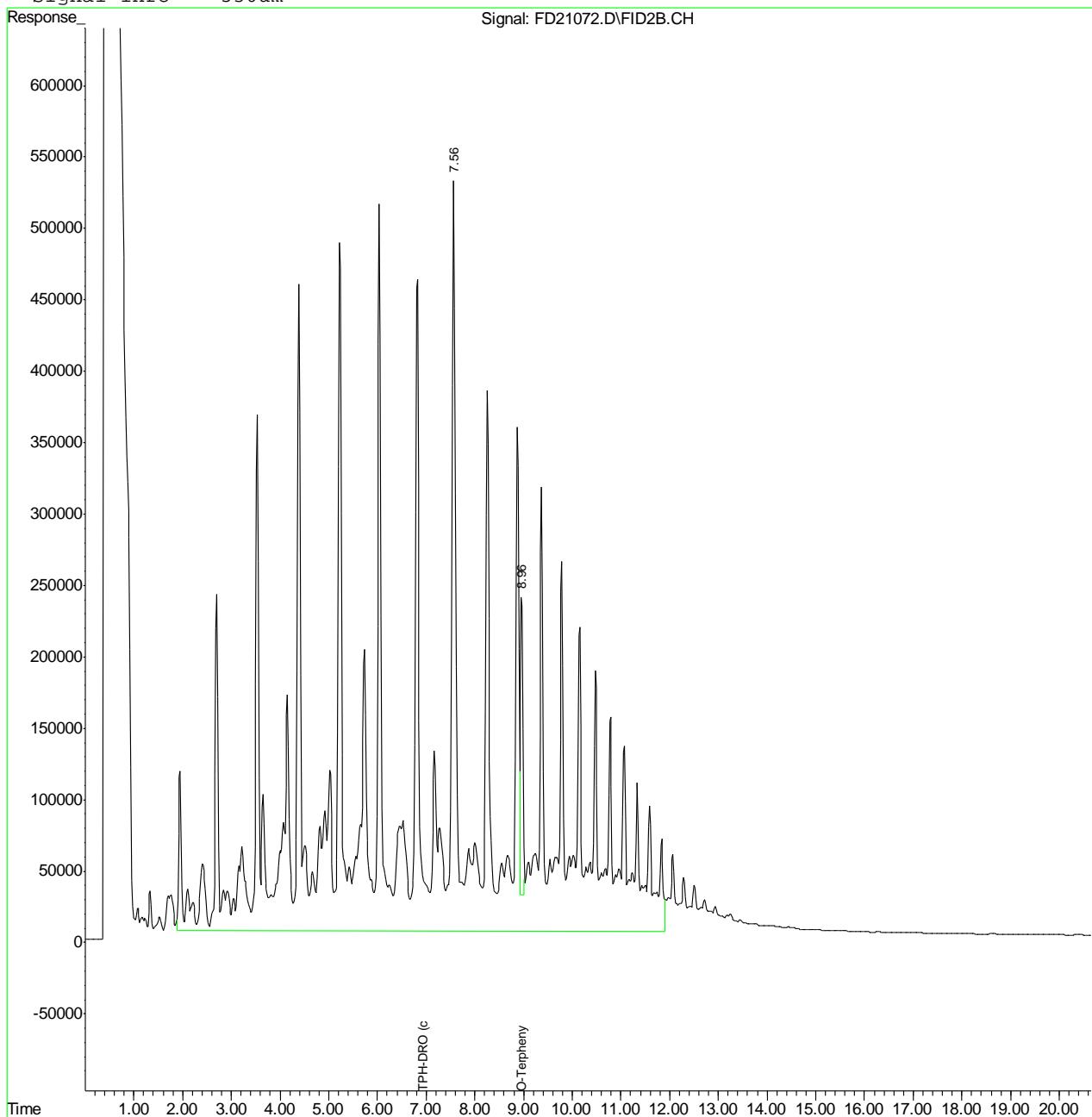
-----  
 (f)=RT Delta > 1/2 Window (m)=manual int.  
 FD21072.D DRO-GFD983R.M Fri Jan 11 08:49:40 2013 GC

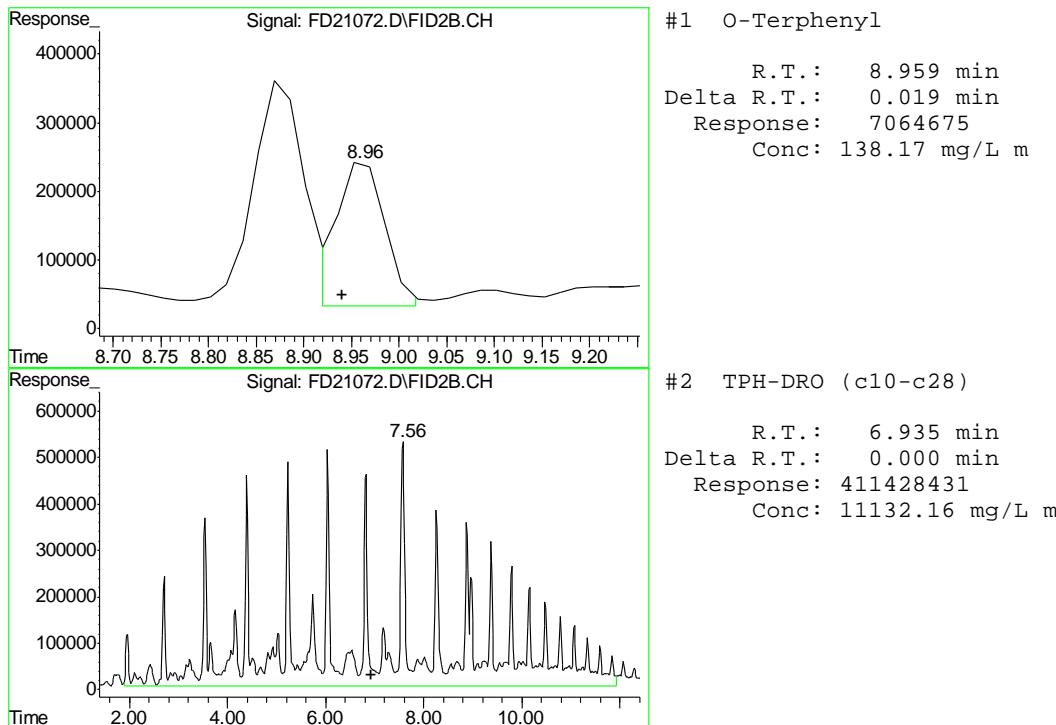
## Quantitation Report (QT Reviewed)

Data File : C:\MSDCHEM\2\DATA\2013\JAN\FD011013.SEC\FD21072.D Vial: 72  
 Acq On : 1-10-2013 09:05:02 PM Operator: ashleyv  
 Sample : D42316-1, 10x Inst : FID5  
 Misc : OP7201,GFD1060,30.06,,,1,10 Multiplr: 1.00  
 IntFile : autoint1.e  
 Quant Time: Jan 11 8:37 2013 Quant Results File: DRO-GFD983R.RES

Quant Method : C:\MSDCHEM\2...\DRO-GFD983R.M (Chemstation Integrator)  
 Title : 8015B TEH  
 Last Update : Tue Jan 08 09:15:06 2013  
 Response via : Multiple Level Calibration  
 DataAcq Meth : DRODUAL.M

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





## Quantitation Report (QT Reviewed)

Data File : C:\MSDCHEM\2\DATA\2013\JAN\FD010913.SEC\FD21006.D Vial: 53  
 Acq On : 1-9-2013 05:11:42 PM Operator: ashleyv  
 Sample : OP7201-MB Inst : FID5  
 Misc : OP7201,GFD1058,30.00,,,1,1 Multiplr: 1.00  
 IntFile : autoint1.e  
 Quant Time: Jan 10 11:11:54 2013 Quant Results File: DRO-GFD983R.RES

Quant Method : C:\MSDCHEM\2...\DRO-GFD983R.M (Chemstation Integrator)  
 Title : 8015B TEH  
 Last Update : Tue Jan 08 09:15:06 2013  
 Response via : Initial Calibration  
 DataAcq Meth : DRODUAL.M

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

Compound	R.T.	Response	Conc Units
<hr/>			
System Monitoring Compounds			
1) S O-Terphenyl	8.98	85813219	1678.342 mg/L
<hr/>			
Target Compounds			
2) H TPH-DRO (c10-c28)	6.93	1255784	33.978 mg/L

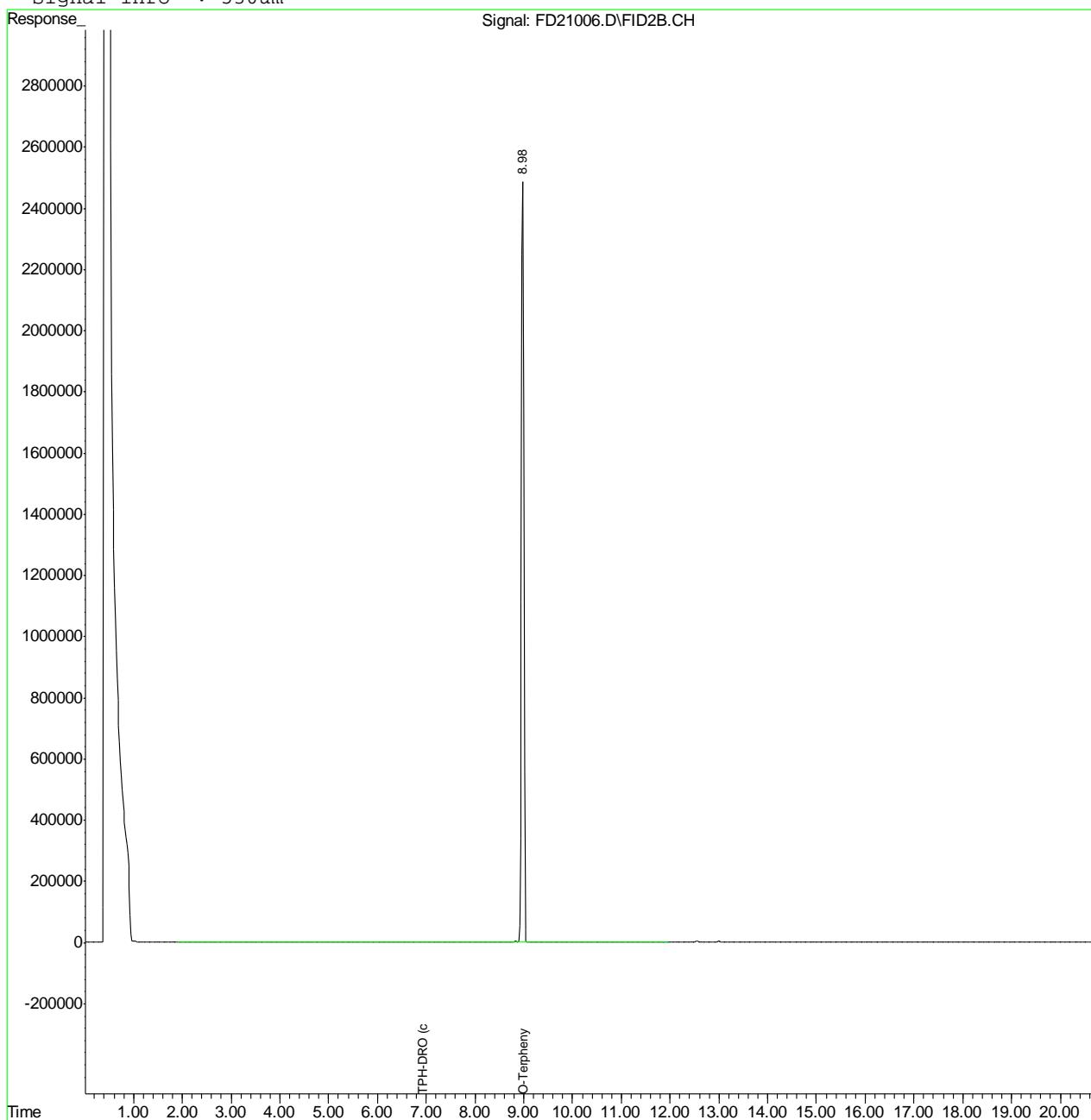
(f)=RT Delta > 1/2 Window (m)=manual int.  
 FD21006.D DRO-GFD983R.M Thu Jan 10 11:19:30 2013 GC

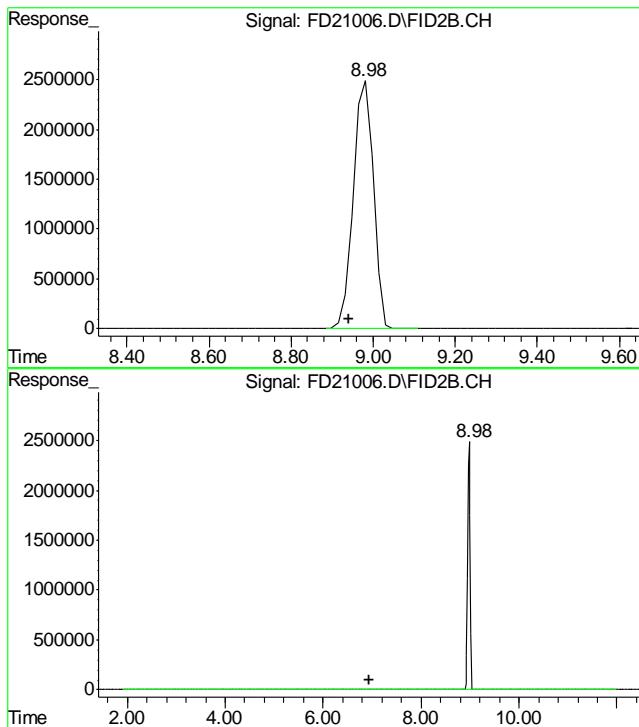
## Quantitation Report (QT Reviewed)

Data File : C:\MSDCHEM\2\DATA\2013\JAN\FD010913.SEC\FD21006.D Vial: 53  
 Acq On : 1-9-2013 05:11:42 PM Operator: ashleyv  
 Sample : OP7201-MB Inst : FID5  
 Misc : OP7201,GFD1058,30.00,,,1,1 Multiplr: 1.00  
 IntFile : autoint1.e  
 Quant Time: Jan 10 11:11 2013 Quant Results File: DRO-GFD983R.RES

Quant Method : C:\MSDCHEM\2...\DRO-GFD983R.M (Chemstation Integrator)  
 Title : 8015B TEH  
 Last Update : Tue Jan 08 09:15:06 2013  
 Response via : Multiple Level Calibration  
 DataAcq Meth : DRODUAL.M

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





#1 O-Terphenyl  
 R.T.: 8.983 min  
 Delta R.T.: 0.043 min  
 Response: 85813219  
 Conc: 1678.34 mg/L

#2 TPH-DRO (c10-c28)  
 R.T.: 6.935 min  
 Delta R.T.: 0.000 min  
 Response: 1255784  
 Conc: 33.98 mg/L



## Metals Analysis

---

### 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: D42316  
Account: XTOKRWR - XTO Energy  
Project: XTO Love Ranch 8

QC Batch ID: MP9199  
Matrix Type: SOLID

Methods: SW846 6010C  
Units: mg/kg

Prep Date:

01/07/13

Metal	RL	IDL	MDL	MB raw	final
Aluminum	10	.96	.57		
Antimony	3.0	.17	.12		
Arsenic	2.5	.44	.56		
Barium	1.0	.01	.11	0.11	<1.0
Beryllium	1.0	.13	.15		
Boron	5.0	.1	.06		
Cadmium	1.0	.06	.036	0.030	<1.0
Calcium	40	.54	9		
Chromium	1.0	.03	.03	0.010	<1.0
Cobalt	0.50	.04	.07		
Copper	1.0	.12	.15	0.0	<1.0
Iron	7.0	.12	.87		
Lead	5.0	.19	.24	-0.30	<5.0
Lithium	0.20	.05	.054		
Magnesium	20	.65	.98		
Manganese	0.50	.12	.022		
Molybdenum	1.0	.21	.08		
Nickel	3.0	.05	.026	-0.050	<3.0
Phosphorus	10	1.4	1.9		
Potassium	200	6.1	7		
Selenium	5.0	.48	.36	0.56	<5.0
Silicon	5.0	.29	.37		
Silver	3.0	.04	.06	0.020	<3.0
Sodium	40	.59	1.9		
Strontium	5.0	.004	.017		
Thallium	1.0	.29	.53		
Tin	5.0	1.2	2		
Titanium	1.0	.01	.038		
Uranium	5.0	.22	.26		
Vanadium	1.0	.02	.036		
Zinc	3.0	.05	.37	0.080	<3.0

Associated samples MP9199: D42316-1

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

BLANK RESULTS SUMMARY  
Part 2 - Method Blanks

Login Number: D42316  
Account: XTOKWR - XTO Energy  
Project: XTO Love Ranch 8

QC Batch ID: MP9199  
Matrix Type: SOLID

Methods: SW846 6010C  
Units: mg/kg

Prep Date:

Metal

(anr) Analyte not requested

## MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D42316  
 Account: XTOKRWR - XTO Energy  
 Project: XTO Love Ranch 8

QC Batch ID: MP9199  
 Matrix Type: SOLID

Methods: SW846 6010C  
 Units: mg/kg

Prep Date:

01/07/13

Metal	D42292-1 Original MS	Spikelot ICPALL2	% Rec	QC Limits
Aluminum				
Antimony				
Arsenic	anr			
Barium	105	291	206	90.1 75-125
Beryllium				
Boron				
Cadmium	0.071	45.3	51.6	87.7 75-125
Calcium				
Chromium	3.4	50.9	51.6	92.1 75-125
Cobalt				
Copper	3.1	49.0	51.6	89.0 75-125
Iron				
Lead	3.2	96.0	103	89.9 75-125
Lithium				
Magnesium				
Manganese				
Molybdenum				
Nickel	5.1	49.9	51.6	86.8 75-125
Phosphorus				
Potassium				
Selenium	0.0	88.0	103	85.3 75-125
Silicon				
Silver	0.0	20.1	20.6	97.4 75-125
Sodium				
Strontium				
Thallium				
Tin				
Titanium				
Uranium				
Vanadium				
Zinc	10.3	54.1	51.6	84.9 75-125

Associated samples MP9199: D42316-1

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

MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D42316  
Account: XTOKWR - XTO Energy  
Project: XTO Love Ranch 8

QC Batch ID: MP9199  
Matrix Type: SOLID

Methods: SW846 6010C  
Units: mg/kg

Prep Date:

Metal

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

## MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D42316  
 Account: XTOKRWR - XTO Energy  
 Project: XTO Love Ranch 8

QC Batch ID: MP9199  
 Matrix Type: SOLID

Methods: SW846 6010C  
 Units: mg/kg

Prep Date:

01/07/13

Metal	D42292-1 Original	MSD	Spikelot ICPALL2	% Rec	MSD RPD	QC Limit
Aluminum						
Antimony						
Arsenic	anr					
Barium	105	291	198	93.8	0.0	20
Beryllium						
Boron						
Cadmium	0.071	43.3	49.6	87.2	4.5	20
Calcium						
Chromium	3.4	48.9	49.6	91.8	4.0	20
Cobalt						
Copper	3.1	47.9	49.6	90.3	2.3	20
Iron						
Lead	3.2	92.2	99.2	89.7	4.0	20
Lithium						
Magnesium						
Manganese						
Molybdenum						
Nickel	5.1	48.1	49.6	86.7	3.7	20
Phosphorus						
Potassium						
Selenium	0.0	84.1	99.2	84.8	4.5	20
Silicon						
Silver	0.0	19.0	19.8	95.8	5.6	20
Sodium						
Strontium						
Thallium						
Tin						
Titanium						
Uranium						
Vanadium						
Zinc	10.3	53.0	49.6	86.1	2.1	20

Associated samples MP9199: D42316-1

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

MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D42316  
Account: XTOKWR - XTO Energy  
Project: XTO Love Ranch 8

QC Batch ID: MP9199  
Matrix Type: SOLID

Methods: SW846 6010C  
Units: mg/kg

Prep Date:

Metal

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

## SPIKE BLANK AND LAB CONTROL SAMPLE SUMMARY

Login Number: D42316  
 Account: XTOKRWR - XTO Energy  
 Project: XTO Love Ranch 8

QC Batch ID: MP9199  
 Matrix Type: SOLID

Methods: SW846 6010C  
 Units: mg/kg

Prep Date:

01/07/13

Metal	BSP Result	Spikelot ICPALL2	% Rec	QC Limits
Aluminum				
Antimony				
Arsenic	anr			
Barium	182	200	91.0	80-120
Beryllium				
Boron				
Cadmium	45.4	50	90.8	80-120
Calcium				
Chromium	48.0	50	96.0	80-120
Cobalt				
Copper	43.8	50	87.6	80-120
Iron				
Lead	93.7	100	93.7	80-120
Lithium				
Magnesium				
Manganese				
Molybdenum				
Nickel	46.1	50	92.2	80-120
Phosphorus				
Potassium				
Selenium	91.2	100	91.2	80-120
Silicon				
Silver	19.7	20	98.5	80-120
Sodium				
Strontium				
Thallium				
Tin				
Titanium				
Uranium				
Vanadium				
Zinc	45.2	50	90.4	80-120

Associated samples MP9199: D42316-1

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

SPIKE BLANK AND LAB CONTROL SAMPLE SUMMARY

Login Number: D42316  
Account: XTOKRWR - XTO Energy  
Project: XTO Love Ranch 8

QC Batch ID: MP9199  
Matrix Type: SOLID

Methods: SW846 6010C  
Units: mg/kg

Prep Date:

Metal

(anr) Analyte not requested

14.1.3  
**14**

## SERIAL DILUTION RESULTS SUMMARY

Login Number: D42316  
 Account: XTOKWR - XTO Energy  
 Project: XTO Love Ranch 8

QC Batch ID: MP9199  
 Matrix Type: SOLID

Methods: SW846 6010C  
 Units: ug/l

Prep Date: 01/07/13

Metal	D42292-1 Original	SDL 1:5	%DIF	QC Limits
Aluminum				
Antimony				
Arsenic	anr			
Barium	1040	1150	11.1*(a)	0-10
Beryllium				
Boron				
Cadmium	0.700	0.00	100.0(b)	0-10
Calcium				
Chromium	33.4	34.0	1.8	0-10
Cobalt				
Copper	30.9	27.5	11.0 (b)	0-10
Iron				
Lead	31.9	10.5	67.1 (b)	0-10
Lithium				
Magnesium				
Manganese				
Molybdenum				
Nickel	50.4	53.0	5.2	0-10
Phosphorus				
Potassium				
Selenium	0.00	0.00	NC	0-10
Silicon				
Silver	0.00	2.00	NC	0-10
Sodium				
Strontium				
Thallium				
Tin				
Titanium				
Uranium				
Vanadium				
Zinc	102	118	15.3*(a)	0-10

Associated samples MP9199: D42316-1

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

SERIAL DILUTION RESULTS SUMMARY

Login Number: D42316  
Account: XTOKRWR - XTO Energy  
Project: XTO Love Ranch 8

QC Batch ID: MP9199  
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: D42316  
Account: XTOKWR - XTO Energy  
Project: XTO Love Ranch 8

QC Batch ID: MP9200  
Matrix Type: SOLID

Methods: SW846 6020A  
Units: mg/kg

Prep Date:

01/07/13

Metal	RL	IDL	MDL	MB raw	final
Aluminum	25	.22	.31		
Antimony	0.20	.0018	.0075		
Arsenic	0.10	.006	.06	0.011	<0.10
Barium	1.0	.0065	.037		
Beryllium	0.10	.016	.09		
Boron	20	1.2	1.2		
Cadmium	0.050	.014	.021		
Calcium	200	7.9	8		
Chromium	1.0	.033	.19		
Cobalt	0.10	.0012	.015		
Copper	1.0	.017	.065		
Iron	20	.8	5		
Lead	0.25	.0011	.024		
Magnesium	50	.44	.85		
Manganese	0.50	.0043	.02		
Molybdenum	0.50	.018	.018		
Nickel	1.0	.0049	.011		
Phosphorus	30	1.4	3.6		
Potassium	100	9.8	10		
Selenium	0.20	.029	.14		
Silver	0.050	.0009	.0065		
Sodium	250	1.5	2.3		
Strontium	10	.036	.036		
Thallium	0.10	.00095	.0095		
Tin	5.0	.023	.34		
Titanium	1.0	.044	.1		
Uranium	0.25	.00085	.001		
Vanadium	2.0	.12	.21		
Zinc	5.0	.033	.35		

Associated samples MP9200: D42316-1

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

14.2.1  
14

## MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D42316  
 Account: XTOKWR - XTO Energy  
 Project: XTO Love Ranch 8

QC Batch ID: MP9200  
 Matrix Type: SOLID

Methods: SW846 6020A  
 Units: mg/kg

Prep Date: 01/07/13

Metal	D42292-1 Original MS	Spikelot ICPALL2	% Rec	QC Limits
Aluminum				
Antimony				
Arsenic	1.2	109	103	104.5 75-125
Barium				
Beryllium				
Boron				
Cadmium	anr			
Calcium				
Chromium	anr			
Cobalt				
Copper	anr			
Iron				
Lead	anr			
Magnesium				
Manganese				
Molybdenum	anr			
Nickel	anr			
Phosphorus	anr			
Potassium	anr			
Selenium	anr			
Silver				
Sodium				
Strontium				
Thallium				
Tin				
Titanium				
Uranium				
Vanadium				
Zinc	anr			

Associated samples MP9200: D42316-1

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: D42316  
 Account: XTOKWR - XTO Energy  
 Project: XTO Love Ranch 8

QC Batch ID: MP9200  
 Matrix Type: SOLID

Methods: SW846 6020A  
 Units: mg/kg

Prep Date: 01/07/13

Metal	D42292-1 Original	MSD	Spikelot ICPALL2	% Rec	MSD RPD	QC Limit
Aluminum						
Antimony						
Arsenic	1.2	100	99.2	99.6	8.6	20
Barium						
Beryllium						
Boron						
Cadmium		anr				
Calcium						
Chromium		anr				
Cobalt						
Copper		anr				
Iron						
Lead		anr				
Magnesium						
Manganese						
Molybdenum		anr				
Nickel		anr				
Phosphorus		anr				
Potassium		anr				
Selenium		anr				
Silver						
Sodium						
Strontium						
Thallium						
Tin						
Titanium						
Uranium						
Vanadium						
Zinc		anr				

Associated samples MP9200: D42316-1

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: D42316  
 Account: XTOKWR - XTO Energy  
 Project: XTO Love Ranch 8

QC Batch ID: MP9200  
 Matrix Type: SOLID

Methods: SW846 6020A  
 Units: mg/kg

Prep Date: 01/07/13

Metal	BSP Result	Spikelot ICPALL2	% Rec	QC Limits
Aluminum				
Antimony				
Arsenic	103	100	103.0	80-120
Barium				
Beryllium				
Boron				
Cadmium	anr			
Calcium				
Chromium	anr			
Cobalt				
Copper	anr			
Iron				
Lead	anr			
Magnesium				
Manganese				
Molybdenum	anr			
Nickel	anr			
Phosphorus	anr			
Potassium	anr			
Selenium	anr			
Silver				
Sodium				
Strontium				
Thallium				
Tin				
Titanium				
Uranium				
Vanadium				
Zinc	anr			

Associated samples MP9200: D42316-1

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

## SERIAL DILUTION RESULTS SUMMARY

Login Number: D42316  
 Account: XTOKWR - XTO Energy  
 Project: XTO Love Ranch 8

QC Batch ID: MP9200  
 Matrix Type: SOLID

Methods: SW846 6020A  
 Units: ug/l

Prep Date: 01/07/13

Metal	D42292-1 Original	SDL 5:25	%DIF	QC Limits
Aluminum				
Antimony				
Arsenic	11.5	10.7	6.8	0-10
Barium				
Beryllium				
Boron				
Cadmium	anr			
Calcium				
Chromium	anr			
Cobalt				
Copper	anr			
Iron				
Lead	anr			
Magnesium				
Manganese				
Molybdenum	anr			
Nickel	anr			
Phosphorus	anr			
Potassium	anr			
Selenium	anr			
Silver				
Sodium				
Strontium				
Thallium				
Tin				
Titanium				
Uranium				
Vanadium				
Zinc	anr			

Associated samples MP9200: D42316-1

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

14.2.4  
**14**

BLANK RESULTS SUMMARY  
Part 2 - Method Blanks

Login Number: D42316  
Account: XTOKWR - XTO Energy  
Project: XTO Love Ranch 8

QC Batch ID: MP9202  
Matrix Type: SOLID

Methods: SW846 7471B  
Units: mg/kg

Prep Date:

01/08/13

Metal	RL	IDL	MDL	MB raw	final
Mercury	0.10	.0011	.0009	-0.00031	<0.10

Associated samples MP9202: D42316-1

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: D42316  
Account: XTOKRWR - XTO Energy  
Project: XTO Love Ranch 8

QC Batch ID: MP9202  
Matrix Type: SOLID

Methods: SW846 7471B  
Units: mg/kg

Prep Date: 01/08/13

Metal	D42292-1 Original MS	Spikelot HGWSR1	QC % Rec	QC Limits
Mercury	0.0062	0.36	0.335	105.6 75-125

Associated samples MP9202: D42316-1

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: D42316  
Account: XTOKWR - XTO Energy  
Project: XTO Love Ranch 8

QC Batch ID: MP9202  
Matrix Type: SOLID

Methods: SW846 7471B  
Units: mg/kg

Prep Date:

01/08/13

Metal	D42292-1 Original	MSD	Spikelot HGWSR1	MSD % Rec	QC RPD	QC Limit
Mercury	0.0062	0.33	0.314	103.0	8.7	20

Associated samples MP9202: D42316-1

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: D42316  
Account: XTOKRWR - XTO Energy  
Project: XTO Love Ranch 8

QC Batch ID: MP9202  
Matrix Type: SOLID

Methods: SW846 7471B  
Units: mg/kg

Prep Date: 01/08/13

Metal	BSP Result	Spikelot HGWSR1	QC % Rec	QC Limits
Mercury	0.41	0.4	102.5	80-120

Associated samples MP9202: D42316-1

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

14.3.3  
**14**

BLANK RESULTS SUMMARY  
Part 2 - Method Blanks

Login Number: D42316  
Account: XTOKWR - XTO Energy  
Project: XTO Love Ranch 8

QC Batch ID: MP9206  
Matrix Type: AQUEOUS

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

Prep Date:

01/08/13

Metal	RL	IDL	MDL	MB raw	final
Aluminum	500	48	130		
Antimony	150	8.5	18		
Arsenic	130	22	42		
Barium	50	.5	9		
Beryllium	50	6.5	16		
Boron	250	5	22		
Cadmium	50	3	3		
Calcium	2000	27	80	10.5	<2000
Chromium	50	1.5	2.8		
Cobalt	25	2	2.1		
Copper	50	6	15		
Iron	350	6	100		
Lead	250	9.5	15		
Lithium	10	2.5			
Magnesium	1000	33	110	11.5	<1000
Manganese	25	6	6		
Molybdenum	50	11	11		
Nickel	150	2.5	2.9		
Phosphorus	500	70	300		
Potassium	5000	310	750		
Selenium	250	24	55		
Silicon	250	15			
Silver	150	2	4.9		
Sodium	2000	30	490	80.5	<2000
Strontium	25	.2	7.5		
Thallium	50	15	43		
Tin	250	60			
Titanium	50	.5			
Uranium	250	11	23		
Vanadium	50	1	2.4		
Zinc	150	2.5	12		

Associated samples MP9206: D42316-1A

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

BLANK RESULTS SUMMARY  
Part 2 - Method Blanks

Login Number: D42316  
Account: XTOKWR - XTO Energy  
Project: XTO Love Ranch 8

QC Batch ID: MP9206  
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: D42316  
 Account: XTOKRWR - XTO Energy  
 Project: XTO Love Ranch 8

QC Batch ID: MP9206  
 Matrix Type: AQUEOUS

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

Prep Date: 01/08/13

Metal	D42292-1A Original MS	Spikelot ICPALL2	% Rec	QC Limits
Aluminum				
Antimony				
Arsenic				
Barium				
Beryllium				
Boron				
Cadmium				
Calcium	243000	362000	125000	95.2
Chromium				
Cobalt				
Copper				
Iron				
Lead				
Lithium				
Magnesium	57300	184000	125000	101.4
Manganese				
Molybdenum				
Nickel				
Phosphorus				
Potassium				
Selenium				
Silicon				
Silver				
Sodium	376000	485000	125000	87.2
Strontium				
Thallium				
Tin				
Titanium				
Uranium				
Vanadium				
Zinc				

Associated samples MP9206: D42316-1A

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

MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D42316  
Account: XTOKWR - XTO Energy  
Project: XTO Love Ranch 8

QC Batch ID: MP9206  
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: D42316  
 Account: XTOKRWR - XTO Energy  
 Project: XTO Love Ranch 8

QC Batch ID: MP9206  
 Matrix Type: AQUEOUS

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

Prep Date: 01/08/13

Metal	D42292-1A Original MSD	Spikelot ICPALL2	MSD % Rec	MSD RPD	QC Limit
Aluminum					
Antimony					
Arsenic					
Barium					
Beryllium					
Boron					
Cadmium					
Calcium	243000	366000	125000	98.4	1.1
Chromium					
Cobalt					
Copper					
Iron					
Lead					
Lithium					
Magnesium	57300	184000	125000	101.4	0.0
Manganese					
Molybdenum					
Nickel					
Phosphorus					
Potassium					
Selenium					
Silicon					
Silver					
Sodium	376000	486000	125000	88.0	0.2
Strontium					
Thallium					
Tin					
Titanium					
Uranium					
Vanadium					
Zinc					

Associated samples MP9206: D42316-1A

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

MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D42316  
Account: XTOKWR - XTO Energy  
Project: XTO Love Ranch 8

QC Batch ID: MP9206  
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: D42316  
 Account: XTOKRWR - XTO Energy  
 Project: XTO Love Ranch 8

QC Batch ID: MP9206  
 Matrix Type: AQUEOUS

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

Prep Date: 01/08/13

Metal	BSP Result	Spikelot ICPALL2	% Rec	QC Limits
Aluminum				
Antimony				
Arsenic				
Barium				
Beryllium				
Boron				
Cadmium				
Calcium	133000	125000	106.4	80-120
Chromium				
Cobalt				
Copper				
Iron				
Lead				
Lithium				
Magnesium	128000	125000	102.4	80-120
Manganese				
Molybdenum				
Nickel				
Phosphorus				
Potassium				
Selenium				
Silicon				
Silver				
Sodium	127000	125000	101.6	80-120
Strontium				
Thallium				
Tin				
Titanium				
Uranium				
Vanadium				
Zinc				

Associated samples MP9206: D42316-1A

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

SPIKE BLANK AND LAB CONTROL SAMPLE SUMMARY

Login Number: D42316  
Account: XTOKWR - XTO Energy  
Project: XTO Love Ranch 8

QC Batch ID: MP9206  
Matrix Type: AQUEOUS

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

Prep Date:

Metal

(anr) Analyte not requested

14.4.3  
**14**

## SERIAL DILUTION RESULTS SUMMARY

Login Number: D42316  
 Account: XTOKWR - XTO Energy  
 Project: XTO Love Ranch 8

QC Batch ID: MP9206  
 Matrix Type: AQUEOUS

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

Prep Date: 01/08/13

Metal	D42292-1A	Original	SDL 1:5	%DIF	QC Limits
Aluminum					
Antimony					
Arsenic					
Barium					
Beryllium					
Boron					
Cadmium					
Calcium	48500	49200		1.3	0-10
Chromium					
Cobalt					
Copper					
Iron					
Lead					
Lithium					
Magnesium	11500	11500		0.3	0-10
Manganese					
Molybdenum					
Nickel					
Phosphorus					
Potassium					
Selenium					
Silicon					
Silver					
Sodium	75300	75900		0.9	0-10
Strontium					
Thallium					
Tin					
Titanium					
Uranium					
Vanadium					
Zinc					

Associated samples MP9206: D42316-1A

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

SERIAL DILUTION RESULTS SUMMARY

Login Number: D42316  
Account: XTOKRWR - XTO Energy  
Project: XTO Love Ranch 8

QC Batch ID: MP9206  
Matrix Type: AQUEOUS

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

Prep Date:

Metal

(anr) Analyte not requested



## General Chemistry

---

### QC Data Summaries

---

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: D42316  
Account: XTOKWR - XTO Energy  
Project: XTO Love Ranch 8

Analyte	Batch ID	RL	MB Result	Units	Spike Amount	BSP Result	BSP %Recov	QC Limits
Chromium, Hexavalent	GP9044/GN18328	1.0	0.0	mg/kg	176.0	160	90.7	80-120%
Specific Conductivity	GP9061/GN18362			umhos/cm	9992	9840	98.5	90-110%
pH	GN18331			su	8.00	7.99	99.9	99.3-100.7%

Associated Samples:

Batch GP9044: D42316-1

Batch GP9061: D42316-1

Batch GN18331: D42316-1

(\*) Outside of QC limits

DUPLICATE RESULTS SUMMARY  
GENERAL CHEMISTRY

Login Number: D42316  
Account: XTOKWR - XTO Energy  
Project: XTO Love Ranch 8

Analyte	Batch ID	QC Sample	Units	Original Result	DUP Result	RPD	QC Limits
Chromium, Hexavalent Redox Potential Vs H2	GP9044/GN18328 GN18332	D42316-1 D42316-1	mg/kg mv	0.0 36.3	0.0 37.5	0.0 3.2	0-20% 0-20%

Associated Samples:  
Batch GP9044: D42316-1  
Batch GN18332: D42316-1  
(\*) Outside of QC limits

MATRIX SPIKE RESULTS SUMMARY  
GENERAL CHEMISTRY

Login Number: D42316  
Account: XTOKWR - XTO Energy  
Project: XTO Love Ranch 8

Analyte	Batch ID	QC Sample	Units	Original Result	Spike Amount	MS Result	%Rec	QC Limits
Chromium, Hexavalent	GP9044/GN18328	D42316-1	mg/kg	0.0	40.0	38.7	96.8	75-125%

Associated Samples:

Batch GP9044: D42316-1

(\*) Outside of QC limits

(N) Matrix Spike Rec. outside of QC limits

MATRIX SPIKE DUPLICATE RESULTS SUMMARY  
GENERAL CHEMISTRY

Login Number: D42316  
Account: XTOKWR - XTO Energy  
Project: XTO Love Ranch 8

Analyte	Batch ID	QC Sample	Units	Original Result	Spike Amount	MSD Result	RPD	QC Limit
Chromium, Hexavalent	GP9044/GN18328	D42316-1	mg/kg	0.0	40.0	39.2	1.2	20%

Associated Samples:

Batch GP9044: D42316-1

(\*) Outside of QC limits

(N) Matrix Spike Rec. outside of QC limits

15.4

15