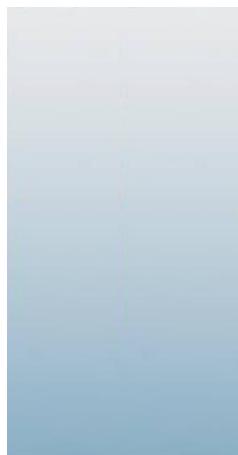




01/17/13



## Technical Report for

**XTO Energy**

**PCU 296-5A**

**1210-04**

**Accutest Job Number: D42556**

**Sampling Date: 01/08/13**

### 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: 193**



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

A handwritten signature in black ink.

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

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

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

This report shall not be reproduced, except in its entirety, without the written approval of Accutest Laboratories.  
Test results relate only to samples analyzed.

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## Sample Summary

XTO Energy

Job No: D42556

PCU 296-5A

Project No: 1210-04

Sample Number	Collected Date	Time By	Received	Matrix Code	Type	Client Sample ID
D42556-1	01/08/13	13:00 DS	01/11/13	SO	Soil	CUT 1 OVERBURDEN
D42556-1A	01/08/13	13:00 DS	01/11/13	SO	Soil	CUT 1 OVERBURDEN
D42556-2	01/08/13	13:10 DS	01/11/13	SO	Soil	CUT 1 CONTENTS
D42556-2A	01/08/13	13:10 DS	01/11/13	SO	Soil	CUT 1 CONTENTS

---

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



## CASE NARRATIVE / CONFORMANCE SUMMARY

**Client:** XTO Energy

**Job No** D42556

**Site:** PCU 296-5A

**Report Date** 1/17/2013 3:09:26 PM

On 01/11/2013, 2 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 D42556 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> V3V1327
------------------	--------------------------

- All samples were analyzed within the recommended method holding time.
- Sample(s) D42512-1MS, D42512-1MSD were used as the QC samples indicated.
- All method blanks for this batch meet method specific criteria.

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

<b>Matrix</b> SO	<b>Batch ID:</b> OP7223
------------------	-------------------------

- All samples were extracted and analyzed within the recommended method holding time.
- Sample(s) D42510-1MS, D42510-1MSD were used as the QC samples indicated.
- All method blanks for this batch meet method specific criteria.
- The matrix spike (MS) recovery(s) of 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 Naphthalene are outside control limits for sample OP7223-MSD. Variability of recovery may be due to sample matrix/homogeneity.

### Volatiles by GC By Method SW846 8015B

<b>Matrix</b> SO	<b>Batch ID:</b> GGB1045
------------------	--------------------------

- All samples were analyzed within the recommended method holding time.
- Sample(s) D42466-11MS, D42466-11MSD were used as the QC samples indicated.
- All method blanks for this batch meet method specific criteria.

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

<b>Matrix</b> SO	<b>Batch ID:</b> OP7232
------------------	-------------------------

- All samples were extracted and analyzed within the recommended method holding time.
- Sample(s) D42562-1MS, D42562-1MSD were used as the QC samples indicated.
- All method blanks for this batch meet method specific criteria.

## Metals By Method SW846 6010C

**Matrix** AQ

**Batch ID:** MP9251

- All samples were digested and analyzed within the recommended method holding time.
- All method blanks for this batch meet method specific criteria.
- Sample(s) D42556-1AMS, D42556-1AMSD, D42556-1ASDL were used as the QC samples for the metals analysis.

**Matrix** SO

**Batch ID:** MP9242

- All samples were digested and analyzed within the recommended method holding time.
- All method blanks for this batch meet method specific criteria.
- Sample(s) D42510-1MS, D42510-1MSD, D42510-1SDL were used as the QC samples for the metals analysis.
- The matrix spike (MS) recovery(s) of Barium are outside control limits. Spike amount low relative to the sample amount. Refer to lab control or spike blank for recovery information.
- The serial dilution RPD(s) for Chromium, Lead, Zinc are outside control limits for sample MP9242-SD1. Probable cause due to sample homogeneity.
- MP9242-SD1 for Chromium: Serial dilution indicates possible matrix interference.
- MP9242-SD1 for Lead: Serial dilution indicates possible matrix interference.
- MP9242-SD1 for Zinc: Serial dilution indicates possible matrix interference.

## Metals By Method SW846 6020A

**Matrix** SO

**Batch ID:** MP9243

- All samples were digested and analyzed within the recommended method holding time.
- All method blanks for this batch meet method specific criteria.
- Sample(s) D42510-1MS, D42510-1SDL, D42510-1MSD were used as the QC samples for the metals analysis.

## Metals By Method SW846 7471B

**Matrix** SO

**Batch ID:** MP9244

- All samples were digested and analyzed within the recommended method holding time.
- All method blanks for this batch meet method specific criteria.
- Sample(s) D42445-1MS, D42445-1MSD were used as the QC samples for the metals analysis.

## Wet Chemistry By Method ASTM D1498-76M

**Matrix** SO

**Batch ID:** GN18422

- Sample(s) D42337-20DUP were used as the QC samples for the Redox Potential Vs H<sub>2</sub> analysis.

## Wet Chemistry By Method SM 2510B-2011 MOD

**Matrix** SO

**Batch ID:** GP9098

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

## Wet Chemistry By Method SM19 2540B M

**Matrix** SO

**Batch ID:** GN18412

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

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

**Matrix** SO

**Batch ID:** GP9086

- All samples were prepared and analyzed within the recommended method holding time.
- All method blanks for this batch meet method specific criteria.
- Sample(s) D42556-1DUP, D42556-1MS, D42556-1MSD were used as the QC samples for the Chromium, Hexavalent analysis.

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

**Matrix** SO

**Batch ID:** R15681

- The data for SW846 3060A/7196A M meets quality control requirements.
- D42556-2 for Chromium, Trivalent: Calculated as: (Chromium) - (Chromium, Hexavalent)

**Matrix** SO

**Batch ID:** R15684

- The data for SW846 3060A/7196A M meets quality control requirements.
- D42556-1 for Chromium, Trivalent: Calculated as: (Chromium) - (Chromium, Hexavalent)

## Wet Chemistry By Method SW846 9045D

**Matrix** SO

**Batch ID:** GN18424

- The following samples were run outside of holding time for method SW846 9045D: D42556-1

**Matrix** SO

**Batch ID:** GN18426

- The following samples were run outside of holding time for method SW846 9045D: D42556-2

## Wet Chemistry By Method USDA HANDBOOK 60

**Matrix** SO

**Batch ID:** MP9251

- D42556-1A for Sodium Adsorption Ratio: Calculated as:  $(Na \text{ meq/L}) / \sqrt{(Ca \text{ meq/L}) + (Mg \text{ meq/L})/2}$
- D42556-2A for Sodium Adsorption Ratio: Calculated as:  $(Na \text{ meq/L}) / \sqrt{(Ca \text{ meq/L}) + (Mg \text{ meq/L})/2}$

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 2

Job Number: D42556  
Account: XTO Energy  
Project: PCU 296-5A  
Collected: 01/08/13

3

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

### D42556-1 CUT 1 OVERBURDEN

TPH-DRO (C10-C28)	45.6	7.5	4.5	mg/kg	SW846-8015B
Arsenic	4.4	0.11		mg/kg	SW846 6020A
Barium	679	1.1		mg/kg	SW846 6010C
Chromium	32.7	1.1		mg/kg	SW846 6010C
Copper	11.8	1.1		mg/kg	SW846 6010C
Lead	10.1	5.7		mg/kg	SW846 6010C
Nickel	17.0	3.4		mg/kg	SW846 6010C
Zinc	36.1	3.4		mg/kg	SW846 6010C
Specific Conductivity	1050	1.0		umhos/cm	SM 2510B-2011 MOD
Chromium, Trivalent <sup>a</sup>	32.7	2.1		mg/kg	SW846 3060A/7196A M
Redox Potential Vs H2	176			mv	ASTM D1498-76M
pH	9.54			su	SW846 9045D

### D42556-1A CUT 1 OVERBURDEN

Calcium	92.3	2.0		mg/l	SW846 6010C
Magnesium	26.0	1.0		mg/l	SW846 6010C
Sodium	107	2.0		mg/l	SW846 6010C
Sodium Adsorption Ratio <sup>b</sup>	2.53			ratio	USDA HANDBOOK 60

### D42556-2 CUT 1 CONTENTS

Toluene	0.396	0.14	0.072	mg/kg	SW846 8260B
Ethylbenzene	0.179	0.14	0.027	mg/kg	SW846 8260B
Xylene (total)	0.854	0.29	0.14	mg/kg	SW846 8260B
Naphthalene	0.185	0.014	0.013	mg/kg	SW846 8270C BY SIM
Pyrene	0.0270	0.010	0.0053	mg/kg	SW846 8270C BY SIM
TPH-GRO (C6-C10)	24.9	14	7.2	mg/kg	SW846 8015B
TPH-DRO (C10-C28)	604	8.1	4.9	mg/kg	SW846-8015B
Arsenic	10.2	0.12		mg/kg	SW846 6020A
Barium	1640	1.2		mg/kg	SW846 6010C
Chromium	24.1	1.2		mg/kg	SW846 6010C
Copper	24.8	1.2		mg/kg	SW846 6010C
Lead	39.4	6.1		mg/kg	SW846 6010C
Nickel	15.4	3.7		mg/kg	SW846 6010C
Zinc	51.3	3.7		mg/kg	SW846 6010C
Specific Conductivity	8300	1.0		umhos/cm	SM 2510B-2011 MOD
Chromium, Trivalent <sup>a</sup>	24.1	2.2		mg/kg	SW846 3060A/7196A M
Redox Potential Vs H2	135			mv	ASTM D1498-76M
pH	10.66			su	SW846 9045D

## Summary of Hits

Page 2 of 2

Job Number: D42556  
Account: XTO Energy  
Project: PCU 296-5A  
Collected: 01/08/13

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

### D42556-2A CUT 1 CONTENTS

Calcium	43.8	2.0	mg/l	SW846 6010C
Sodium	1440	2.0	mg/l	SW846 6010C
Sodium Adsorption Ratio <sup>b</sup>	59.2		ratio	USDA HANDBOOK 60

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

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



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## Sample Results

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### Report of Analysis

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**Report of Analysis**

Page 1 of 1

**Client Sample ID:** CUT 1 OVERBURDEN**Lab Sample ID:** D42556-1**Matrix:** SO - Soil**Method:** SW846 8260B**Project:** PCU 296-5A**Date Sampled:** 01/08/13**Date Received:** 01/11/13**Percent Solids:** 88.2

	<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	3V22570.D	1	01/12/13	BD	n/a	n/a	V3V1327
Run #2							

	<b>Initial Weight</b>	<b>Final Volume</b>	<b>Methanol Aliquot</b>
Run #1	5.07 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	0.063	0.031	mg/kg	
108-88-3	Toluene	ND	0.13	0.063	mg/kg	
100-41-4	Ethylbenzene	ND	0.13	0.024	mg/kg	
1330-20-7	Xylene (total)	ND	0.25	0.13	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	86%		64-130%
460-00-4	4-Bromofluorobenzene	104%		62-131%
17060-07-0	1,2-Dichloroethane-D4	91%		70-130%

ND = Not detected MDL - Method Detection Limit

J = Indicates an estimated value

RL = Reporting Limit

B = Indicates analyte found in associated method blank

E = Indicates value exceeds calibration range

N = Indicates presumptive evidence of a compound

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**Report of Analysis**

Page 1 of 1

<b>Client Sample ID:</b>	CUT 1 OVERBURDEN	<b>Date Sampled:</b>	01/08/13
<b>Lab Sample ID:</b>	D42556-1	<b>Date Received:</b>	01/11/13
<b>Matrix:</b>	SO - Soil	<b>Percent Solids:</b>	88.2
<b>Method:</b>	SW846 8270C BY SIM	SW846 3546	
<b>Project:</b>	PCU 296-5A		

	<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	3G12980.D	1	01/15/13	DC	01/14/13	OP7223	E3G621
Run #2							

	<b>Initial Weight</b>	<b>Final Volume</b>
Run #1	30.1 g	1.0 ml
Run #2		

**COGCC Table 910-1 PAH List**

<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	0.0094	0.0049	mg/kg	
120-12-7	Anthracene	ND	0.0094	0.0049	mg/kg	
56-55-3	Benzo(a)anthracene	ND	0.0094	0.0049	mg/kg	
205-99-2	Benzo(b)fluoranthene	ND	0.0094	0.0049	mg/kg	
207-08-9	Benzo(k)fluoranthene	ND	0.0094	0.0049	mg/kg	
50-32-8	Benzo(a)pyrene	ND	0.0094	0.0049	mg/kg	
218-01-9	Chrysene	ND	0.0094	0.0049	mg/kg	
53-70-3	Dibenz(a,h)anthracene	ND	0.0094	0.0049	mg/kg	
206-44-0	Fluoranthene	ND	0.0094	0.0049	mg/kg	
86-73-7	Fluorene	ND	0.0094	0.0049	mg/kg	
193-39-5	Indeno(1,2,3-cd)pyrene	ND	0.0094	0.0049	mg/kg	
91-20-3	Naphthalene	ND	0.013	0.012	mg/kg	
129-00-0	Pyrene	ND	0.0094	0.0049	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	83%		10-159%
321-60-8	2-Fluorobiphenyl	87%		19-131%
1718-51-0	Terphenyl-d14	108%		18-150%

ND = Not detected MDL - Method Detection Limit

J = Indicates an estimated value

RL = Reporting Limit

B = Indicates analyte found in associated method blank

E = Indicates value exceeds calibration range

N = Indicates presumptive evidence of a compound

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**Report of Analysis**

Page 1 of 1

**Client Sample ID:** CUT 1 OVERBURDEN**Lab Sample ID:** D42556-1**Matrix:** SO - Soil**Method:** SW846 8015B**Project:** PCU 296-5A**Date Sampled:** 01/08/13**Date Received:** 01/11/13**Percent Solids:** 88.2

	<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	GB19132.D	1	01/11/13	SK	n/a	n/a	GGB1045
Run #2							

	<b>Initial Weight</b>	<b>Final Volume</b>	<b>Methanol Aliquot</b>
Run #1	5.1 g	5.0 ml	100 ul
Run #2			

<b>CAS No.</b>	<b>Compound</b>	<b>Result</b>	<b>RL</b>	<b>MDL</b>	<b>Units</b>	<b>Q</b>
	TPH-GRO (C6-C10)	ND	13	6.3	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	93%		60-140%		

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

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

Accutest Laboratories

**Report of Analysis**

Page 1 of 1

**Client Sample ID:** CUT 1 OVERBURDEN  
**Lab Sample ID:** D42556-1  
**Matrix:** SO - Soil  
**Method:** SW846-8015B SW846 3546  
**Project:** PCU 296-5A

**Date Sampled:** 01/08/13  
**Date Received:** 01/11/13  
**Percent Solids:** 88.2

	<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	FD21152.D	1	01/15/13	AV	01/15/13	OP7232	GFD1064
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)	45.6	7.5	4.5	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	70%		35-130%		

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

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

**Report of Analysis**

Page 1 of 1

**Client Sample ID:** CUT 1 OVERBURDEN**Lab Sample ID:** D42556-1**Matrix:** SO - Soil**Project:** PCU 296-5A**Date Sampled:** 01/08/13**Date Received:** 01/11/13**Percent Solids:** 88.2**Metals Analysis**

Analyte	Result	RL	Units	DF	Prep	Analyzed By	Method	Prep Method
Arsenic	4.4	0.11	mg/kg	5	01/14/13	01/17/13 JB	SW846 6020A <sup>3</sup>	SW846 3050B <sup>5</sup>
Barium	679	1.1	mg/kg	1	01/14/13	01/14/13 JB	SW846 6010C <sup>1</sup>	SW846 3050B <sup>4</sup>
Cadmium	< 1.1	1.1	mg/kg	1	01/14/13	01/14/13 JB	SW846 6010C <sup>1</sup>	SW846 3050B <sup>4</sup>
Chromium	32.7	1.1	mg/kg	1	01/14/13	01/14/13 JB	SW846 6010C <sup>1</sup>	SW846 3050B <sup>4</sup>
Copper	11.8	1.1	mg/kg	1	01/14/13	01/14/13 JB	SW846 6010C <sup>1</sup>	SW846 3050B <sup>4</sup>
Lead	10.1	5.7	mg/kg	1	01/14/13	01/14/13 JB	SW846 6010C <sup>1</sup>	SW846 3050B <sup>4</sup>
Mercury	< 0.094	0.094	mg/kg	1	01/15/13	01/15/13 JB	SW846 7471B <sup>2</sup>	SW846 7471B <sup>6</sup>
Nickel	17.0	3.4	mg/kg	1	01/14/13	01/14/13 JB	SW846 6010C <sup>1</sup>	SW846 3050B <sup>4</sup>
Selenium	< 5.7	5.7	mg/kg	1	01/14/13	01/14/13 JB	SW846 6010C <sup>1</sup>	SW846 3050B <sup>4</sup>
Silver	< 3.4	3.4	mg/kg	1	01/14/13	01/14/13 JB	SW846 6010C <sup>1</sup>	SW846 3050B <sup>4</sup>
Zinc	36.1	3.4	mg/kg	1	01/14/13	01/14/13 JB	SW846 6010C <sup>1</sup>	SW846 3050B <sup>4</sup>

- (1) Instrument QC Batch: MA3169
- (2) Instrument QC Batch: MA3177
- (3) Instrument QC Batch: MA3182
- (4) Prep QC Batch: MP9242
- (5) Prep QC Batch: MP9243
- (6) Prep QC Batch: MP9244

RL = Reporting Limit

**Report of Analysis**

Page 1 of 1

**Client Sample ID:** CUT 1 OVERBURDEN**Lab Sample ID:** D42556-1**Matrix:** SO - Soil**Project:** PCU 296-5A**Date Sampled:** 01/08/13**Date Received:** 01/11/13**Percent Solids:** 88.2**General Chemistry**

Analyte	Result	RL	Units	DF	Analyzed	By	Method
<b>prep: DEPT.OF AG, BOOK N9</b>							
Specific Conductivity	1050	1.0	umhos/cm	1	01/15/13	KB	SM 2510B-2011 MOD
Chromium, Hexavalent	< 1.0	1.0	mg/kg	1	01/14/13	KB	SW846 3060A/7196A
Chromium, Trivalent <sup>a</sup>	32.7	2.1	mg/kg	1	01/14/13 17:24	JB	SW846 3060A/7196A M
Redox Potential Vs H2	176		mv	1	01/14/13	JD	ASTM D1498-76M
Solids, Percent	88.2		%	1	01/14/13	SWT	SM19 2540B M
pH	9.54		su	1	01/14/13 14:20	JD	SW846 9045D

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

RL = Reporting Limit

**Report of Analysis**

Page 1 of 1

**Client Sample ID:** CUT 1 OVERBURDEN**Lab Sample ID:** D42556-1A**Matrix:** SO - Soil**Project:** PCU 296-5A**Date Sampled:** 01/08/13**Date Received:** 01/11/13**Percent Solids:** 88.2**SAR Metals Analysis**

Analyte	Result	RL	Units	DF	Prep	Analyzed By	Method	Prep Method
Calcium	92.3	2.0	mg/l	1	01/15/13	01/16/13 JB	SW846 6010C <sup>1</sup>	SW846 3010A/M <sup>2</sup>
Magnesium	26.0	1.0	mg/l	1	01/15/13	01/16/13 JB	SW846 6010C <sup>1</sup>	SW846 3010A/M <sup>2</sup>
Sodium	107	2.0	mg/l	1	01/15/13	01/16/13 JB	SW846 6010C <sup>1</sup>	SW846 3010A/M <sup>2</sup>

(1) Instrument QC Batch: MA3180

(2) Prep QC Batch: MP9251

RL = Reporting Limit

**Report of Analysis**

Page 1 of 1

**Client Sample ID:** CUT 1 OVERBURDEN**Lab Sample ID:** D42556-1A**Matrix:** SO - Soil**Project:** PCU 296-5A**Date Sampled:** 01/08/13**Date Received:** 01/11/13**Percent Solids:** 88.2**General Chemistry**

Analyte	Result	RL	Units	DF	Analyzed	By	Method
Sodium Adsorption Ratio <sup>a</sup>	2.53		ratio	1	01/16/13 12:27	JB	USDA HANDBOOK 60

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

RL = Reporting Limit

Accutest Laboratories

**Report of Analysis**

Page 1 of 1

4.3  
4**Client Sample ID:** CUT 1 CONTENTS**Lab Sample ID:** D42556-2**Matrix:** SO - Soil**Method:** SW846 8260B**Project:** PCU 296-5A**Date Sampled:** 01/08/13**Date Received:** 01/11/13**Percent Solids:** 81.8

	<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	3V22571.D	1	01/12/13	BD	n/a	n/a	V3V1327
Run #2							

	<b>Initial Weight</b>	<b>Final Volume</b>	<b>Methanol Aliquot</b>
Run #1	5.06 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	0.072	0.036	mg/kg	
108-88-3	Toluene	0.396	0.14	0.072	mg/kg	
100-41-4	Ethylbenzene	0.179	0.14	0.027	mg/kg	
1330-20-7	Xylene (total)	0.854	0.29	0.14	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	87%		64-130%
460-00-4	4-Bromofluorobenzene	105%		62-131%
17060-07-0	1,2-Dichloroethane-D4	87%		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

4.3  
4

<b>Client Sample ID:</b>	CUT 1 CONTENTS	<b>Date Sampled:</b>	01/08/13
<b>Lab Sample ID:</b>	D42556-2	<b>Date Received:</b>	01/11/13
<b>Matrix:</b>	SO - Soil	<b>Percent Solids:</b>	81.8
<b>Method:</b>	SW846 8270C BY SIM	SW846 3546	
<b>Project:</b>	PCU 296-5A		

	<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	3G12987.D	1	01/15/13	DC	01/14/13	OP7223	E3G621
Run #2							

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

**COGCC Table 910-1 PAH List**

<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	0.010	0.0053	mg/kg	
120-12-7	Anthracene	ND	0.010	0.0053	mg/kg	
56-55-3	Benzo(a)anthracene	ND	0.010	0.0053	mg/kg	
205-99-2	Benzo(b)fluoranthene	ND	0.010	0.0053	mg/kg	
207-08-9	Benzo(k)fluoranthene	ND	0.010	0.0053	mg/kg	
50-32-8	Benzo(a)pyrene	ND	0.010	0.0053	mg/kg	
218-01-9	Chrysene	ND	0.010	0.0053	mg/kg	
53-70-3	Dibenz(a,h)anthracene	ND	0.010	0.0053	mg/kg	
206-44-0	Fluoranthene	ND	0.010	0.0053	mg/kg	
86-73-7	Fluorene	ND	0.010	0.0053	mg/kg	
193-39-5	Indeno(1,2,3-cd)pyrene	ND	0.010	0.0053	mg/kg	
91-20-3	Naphthalene	0.185	0.014	0.013	mg/kg	
129-00-0	Pyrene	0.0270	0.010	0.0053	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	58%		10-159%
321-60-8	2-Fluorobiphenyl	82%		19-131%
1718-51-0	Terphenyl-d14	89%		18-150%

ND = Not detected MDL - Method Detection Limit

J = Indicates an estimated value

RL = Reporting Limit

B = Indicates analyte found in associated method blank

E = Indicates value exceeds calibration range

N = Indicates presumptive evidence of a compound

Accutest Laboratories

**Report of Analysis**

Page 1 of 1

4.3  
4**Client Sample ID:** CUT 1 CONTENTS**Lab Sample ID:** D42556-2**Matrix:** SO - Soil**Method:** SW846 8015B**Project:** PCU 296-5A**Date Sampled:** 01/08/13**Date Received:** 01/11/13**Percent Solids:** 81.8

	<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	GB19133.D	1	01/11/13	SK	n/a	n/a	GGB1045
Run #2							

	<b>Initial Weight</b>	<b>Final Volume</b>	<b>Methanol Aliquot</b>
Run #1	5.1 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)	24.9	14	7.2	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	85%			60-140%	

ND = Not detected MDL - Method Detection Limit

J = Indicates an estimated value

RL = Reporting Limit

B = Indicates analyte found in associated method blank

E = Indicates value exceeds calibration range

N = Indicates presumptive evidence of a compound

Accutest Laboratories

**Report of Analysis**

Page 1 of 1

4.3  
4

**Client Sample ID:** CUT 1 CONTENTS  
**Lab Sample ID:** D42556-2  
**Matrix:** SO - Soil  
**Method:** SW846-8015B SW846 3546  
**Project:** PCU 296-5A

**Date Sampled:** 01/08/13  
**Date Received:** 01/11/13  
**Percent Solids:** 81.8

	<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	FD21153.D	1	01/15/13	AV	01/15/13	OP7232	GFD1064
Run #2							

	<b>Initial Weight</b>	<b>Final Volume</b>
Run #1	30.0 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)	604	8.1	4.9	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	81%		35-130%		

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

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

**Report of Analysis**

Page 1 of 1

**Client Sample ID:** CUT 1 CONTENTS**Lab Sample ID:** D42556-2**Matrix:** SO - Soil**Project:** PCU 296-5A**Date Sampled:** 01/08/13**Date Received:** 01/11/13**Percent Solids:** 81.8**Metals Analysis**

Analyte	Result	RL	Units	DF	Prep	Analyzed By	Method	Prep Method
Arsenic	10.2	0.12	mg/kg	5	01/14/13	01/17/13 JB	SW846 6020A <sup>3</sup>	SW846 3050B <sup>5</sup>
Barium	1640	1.2	mg/kg	1	01/14/13	01/14/13 JB	SW846 6010C <sup>1</sup>	SW846 3050B <sup>4</sup>
Cadmium	< 1.2	1.2	mg/kg	1	01/14/13	01/14/13 JB	SW846 6010C <sup>1</sup>	SW846 3050B <sup>4</sup>
Chromium	24.1	1.2	mg/kg	1	01/14/13	01/14/13 JB	SW846 6010C <sup>1</sup>	SW846 3050B <sup>4</sup>
Copper	24.8	1.2	mg/kg	1	01/14/13	01/14/13 JB	SW846 6010C <sup>1</sup>	SW846 3050B <sup>4</sup>
Lead	39.4	6.1	mg/kg	1	01/14/13	01/14/13 JB	SW846 6010C <sup>1</sup>	SW846 3050B <sup>4</sup>
Mercury	< 0.10	0.10	mg/kg	1	01/15/13	01/15/13 JB	SW846 7471B <sup>2</sup>	SW846 7471B <sup>6</sup>
Nickel	15.4	3.7	mg/kg	1	01/14/13	01/14/13 JB	SW846 6010C <sup>1</sup>	SW846 3050B <sup>4</sup>
Selenium	< 6.1	6.1	mg/kg	1	01/14/13	01/14/13 JB	SW846 6010C <sup>1</sup>	SW846 3050B <sup>4</sup>
Silver	< 3.7	3.7	mg/kg	1	01/14/13	01/14/13 JB	SW846 6010C <sup>1</sup>	SW846 3050B <sup>4</sup>
Zinc	51.3	3.7	mg/kg	1	01/14/13	01/14/13 JB	SW846 6010C <sup>1</sup>	SW846 3050B <sup>4</sup>

- (1) Instrument QC Batch: MA3169
- (2) Instrument QC Batch: MA3177
- (3) Instrument QC Batch: MA3182
- (4) Prep QC Batch: MP9242
- (5) Prep QC Batch: MP9243
- (6) Prep QC Batch: MP9244

RL = Reporting Limit

**Report of Analysis**

Page 1 of 1

**Client Sample ID:** CUT 1 CONTENTS**Lab Sample ID:** D42556-2**Matrix:** SO - Soil**Project:** PCU 296-5A**Date Sampled:** 01/08/13**Date Received:** 01/11/13**Percent Solids:** 81.8**General Chemistry**

Analyte	Result	RL	Units	DF	Analyzed	By	Method
<b>prep: DEPT.OF AG, BOOK N9</b>							
Specific Conductivity	8300	1.0	umhos/cm	1	01/15/13	KB	SM 2510B-2011 MOD
Chromium, Hexavalent	< 1.0	1.0	mg/kg	1	01/14/13	KB	SW846 3060A/7196A
Chromium, Trivalent <sup>a</sup>	24.1	2.2	mg/kg	1	01/14/13 17:34	JB	SW846 3060A/7196A M
Redox Potential Vs H2	135		mv	1	01/14/13	JD	ASTM D1498-76M
Solids, Percent	81.8		%	1	01/14/13	SWT	SM19 2540B M
pH	10.66		su	1	01/14/13 14:20	JD	SW846 9045D

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

RL = Reporting Limit

**Report of Analysis**

Page 1 of 1

<b>Client Sample ID:</b>	CUT 1 CONTENTS	<b>Date Sampled:</b>	01/08/13
<b>Lab Sample ID:</b>	D42556-2A	<b>Date Received:</b>	01/11/13
<b>Matrix:</b>	SO - Soil	<b>Percent Solids:</b>	81.8
<b>Project:</b>	PCU 296-5A		

**SAR Metals Analysis**

Analyte	Result	RL	Units	DF	Prep	Analyzed By	Method	Prep Method
Calcium	43.8	2.0	mg/l	1	01/15/13	01/16/13 JB	SW846 6010C <sup>1</sup>	SW846 3010A/M <sup>2</sup>
Magnesium	< 1.0	1.0	mg/l	1	01/15/13	01/16/13 JB	SW846 6010C <sup>1</sup>	SW846 3010A/M <sup>2</sup>
Sodium	1440	2.0	mg/l	1	01/15/13	01/16/13 JB	SW846 6010C <sup>1</sup>	SW846 3010A/M <sup>2</sup>

(1) Instrument QC Batch: MA3180

(2) Prep QC Batch: MP9251

RL = Reporting Limit

**Report of Analysis**

Page 1 of 1

**Client Sample ID:** CUT 1 CONTENTS**Lab Sample ID:** D42556-2A**Matrix:** SO - Soil**Project:** PCU 296-5A**Date Sampled:** 01/08/13**Date Received:** 01/11/13**Percent Solids:** 81.8**General Chemistry**

Analyte	Result	RL	Units	DF	Analyzed	By	Method
Sodium Adsorption Ratio <sup>a</sup>	59.2		ratio	1	01/16/13 13:14	JB	USDA HANDBOOK 60

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

RL = Reporting Limit



## Misc. Forms

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### Custody Documents and Other Forms

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Includes the following where applicable:

- Chain of Custody



## CHAIN OF CUSTODY

PAGE 1 OF 1

4036 Youngfield Street, Wheat Ridge, CO 80033  
TEL. 303-425-6021 FAX: 303-425-6854

D42556: Chain of Custody  
Page 1 of 2



## Accutest Laboratories Sample Receipt Summary

Accutest Job Number: D42556

Client: KRW

Immediate Client Services Action Required: No

Date / Time Received: 1/11/2013 3:30:00 PM

No. Coolers:

1

Client Service Action Required at Login: No

Project: XTO PCU 296-5A

Airbill #'s: HDCO

**Cooler Security**Y or N

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

**Cooler Temperature**Y or N

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

**Quality Control Preservation**Y or N

N/A

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

**Sample Integrity - Documentation**Y or N

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

**Sample Integrity - Condition**Y or N

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

**Sample Integrity - Instructions**Y or N

N/A

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

Comments

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

5.1

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**D42556: Chain of Custody****Page 2 of 2**



## GC/MS Volatiles

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### QC Data Summaries

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Includes the following where applicable:

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



## Method Blank Summary

Page 1 of 1

**Job Number:** D42556  
**Account:** XTOKRWR XTO Energy  
**Project:** PCU 296-5A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
V3V1327-MB	3V22560.D	1	01/12/13	BD	n/a	n/a	V3V1327

**The QC reported here applies to the following samples:**

**Method:** SW846 8260B

D42556-1, D42556-2

6.11

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	89%
460-00-4	4-Bromofluorobenzene	97%
17060-07-0	1,2-Dichloroethane-D4	98%

## Blank Spike Summary

Page 1 of 1

Job Number: D42556

Account: XTOKWR XTO Energy

Project: PCU 296-5A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
V3V1327-BS	3V22561.D	1	01/12/13	BD	n/a	n/a	V3V1327

The QC reported here applies to the following samples:

Method: SW846 8260B

D42556-1, D42556-2

CAS No.	Compound	Spike ug/kg	BSP ug/kg	BSP %	Limits
71-43-2	Benzene	50	47.4	95	70-130
100-41-4	Ethylbenzene	50	48.7	97	70-130
108-88-3	Toluene	50	47.3	95	70-130
1330-20-7	Xylene (total)	150	149	99	70-130

CAS No.	Surrogate Recoveries	BSP	Limits
2037-26-5	Toluene-D8	91%	64-130%
460-00-4	4-Bromofluorobenzene	109%	62-131%
17060-07-0	1,2-Dichloroethane-D4	88%	70-130%

\* = Outside of Control Limits.

# Matrix Spike/Matrix Spike Duplicate Summary

Page 1 of 1

Job Number: D42556

Account: XTOKWR XTO Energy

Project: PCU 296-5A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
D42512-1MS	3V22564.D	1	01/12/13	BD	n/a	n/a	V3V1327
D42512-1MSD	3V22565.D	1	01/12/13	BD	n/a	n/a	V3V1327
D42512-1	3V22563.D	1	01/12/13	BD	n/a	n/a	V3V1327

The QC reported here applies to the following samples:

Method: SW846 8260B

D42556-1, D42556-2

CAS No.	Compound	D42512-1		Spike	MS	MS	MSD	MSD	Limits	
		ug/kg	Q	ug/kg	ug/kg	%	ug/kg	%	RPD	Rec/RPD
71-43-2	Benzene	72.3		3580	3400	93	3340	91	2	64-139/30
100-41-4	Ethylbenzene	35.0	J	3580	3530	98	3500	97	1	68-136/30
108-88-3	Toluene	138	J	3580	3330	89	3300	88	1	60-130/30
1330-20-7	Xylene (total)	212	J	10700	10800	99	10800	99	0	58-142/30

CAS No.	Surrogate Recoveries	MS	MSD	D42512-1	Limits
2037-26-5	Toluene-D8	86%	87%	84%	64-130%
460-00-4	4-Bromofluorobenzene	113%	112%	104%	62-131%
17060-07-0	1,2-Dichloroethane-D4	90%	86%	90%	70-130%

\* = Outside of Control Limits.



GC/MS Volatiles

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Raw Data

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## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\V3011113.S\  
 Data File : 3V22570.D  
 Acq On : 12 Jan 2013 7:13 am  
 Operator : BRETD  
 Sample : D42556-1  
 Misc : MS5218,V3V1327,5.066,,100,5,1  
 ALS Vial : 37 Sample Multiplier: 1

Quant Time: Jan 15 08:28:51 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	11.861	168	340365	50.00	ug/l	0.00
35) 1,4-Difluorobenzene	12.656	114	527966	50.00	ug/l	0.00
53) Chlorobenzene-d5	15.296	117	592540	50.00	ug/l	0.00
74) 1,4-Dichlorobenzene-d4	17.285	152	351094	50.00	ug/l	0.00

System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	12.245	102	34738	45.60	ug/l	0.00
Spiked Amount 50.000	Range 70 - 130		Recovery	=	91.20%	
61) Toluene-d8	14.051	98	612876	42.96	ug/l	0.00
Spiked Amount 50.000	Range 70 - 130		Recovery	=	85.92%	
69) 4-Bromofluorobenzene	16.246	95	320780	52.20	ug/l	0.00
Spiked Amount 50.000	Range 70 - 130		Recovery	=	104.40%	

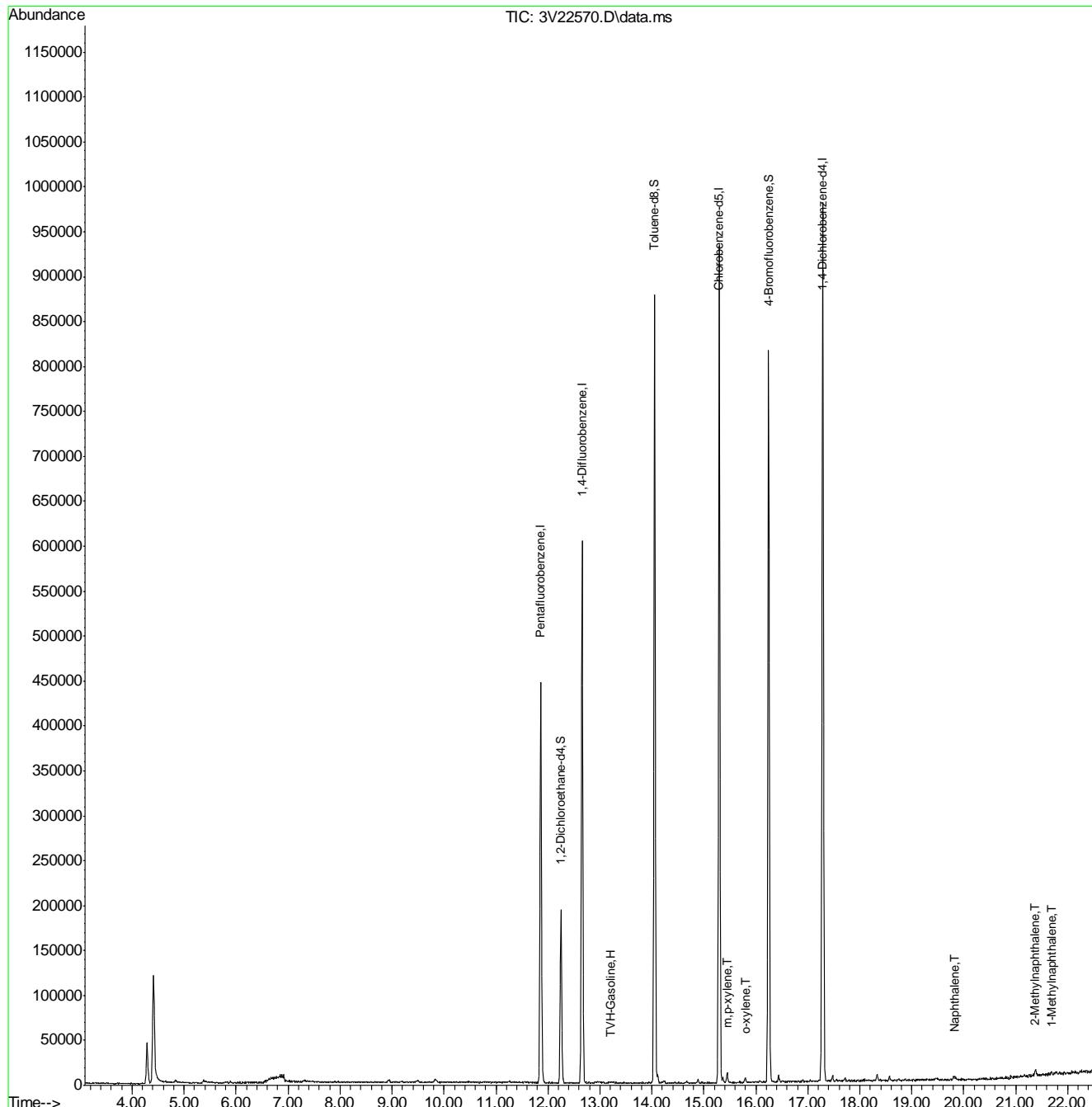
Target Compounds					Qvalue
1) TVH-Gasoline	13.200	TIC	63343m	2.16	ug/l
72) m,p-xylene	15.453	106	4231	0.47	ug/l 92
73) o-xylene	15.796	106	1803	0.20	ug/l 94
91) Naphthalene	19.841	128	3513	1.21	ug/l 100
94) 2-Methylnaphthalene	21.381	142	4567	0.56	ug/l 91
95) 1-Methylnaphthalene	21.689	142	2164	0.28	ug/l 93

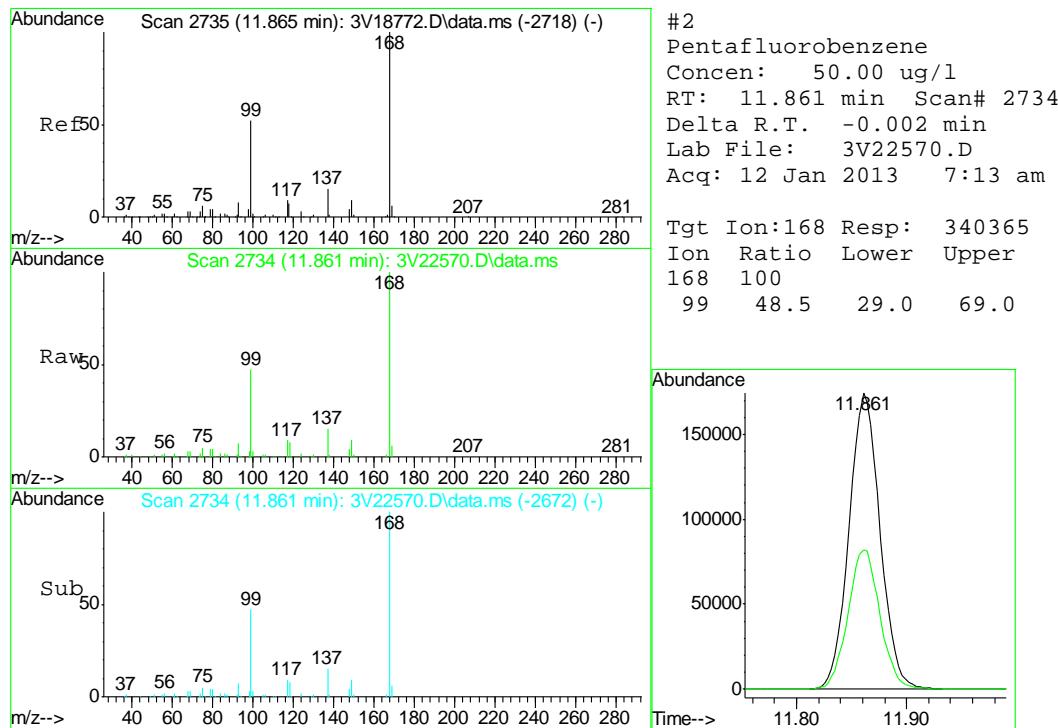
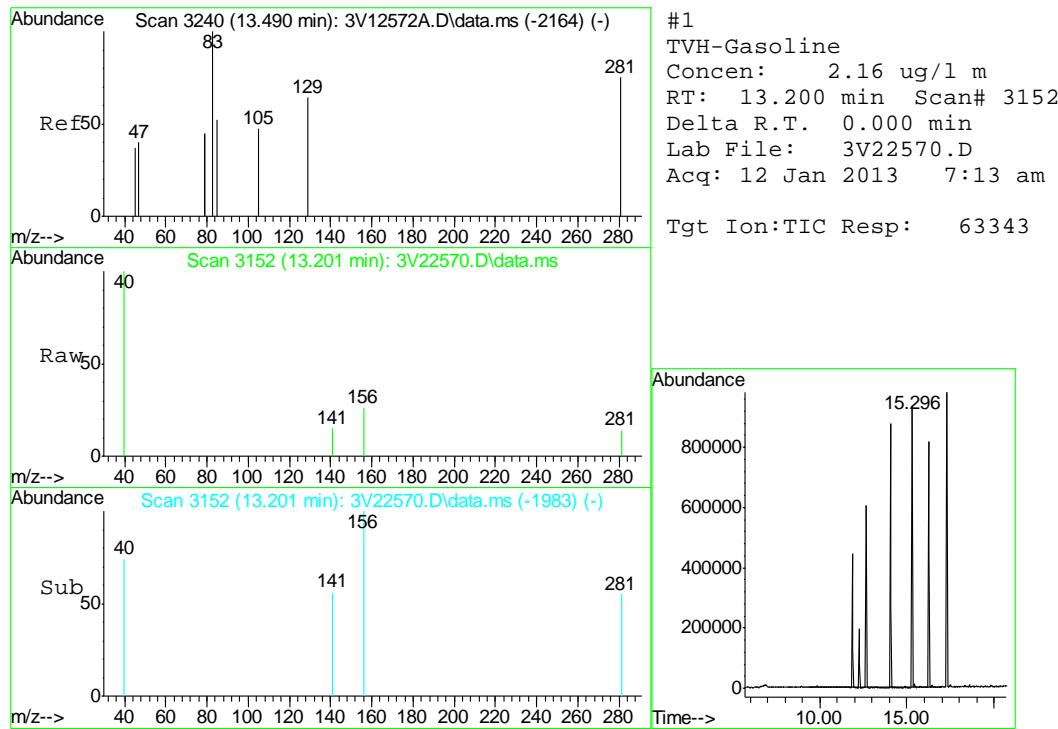
(#) = qualifier out of range (m) = manual integration (+) = signals summed

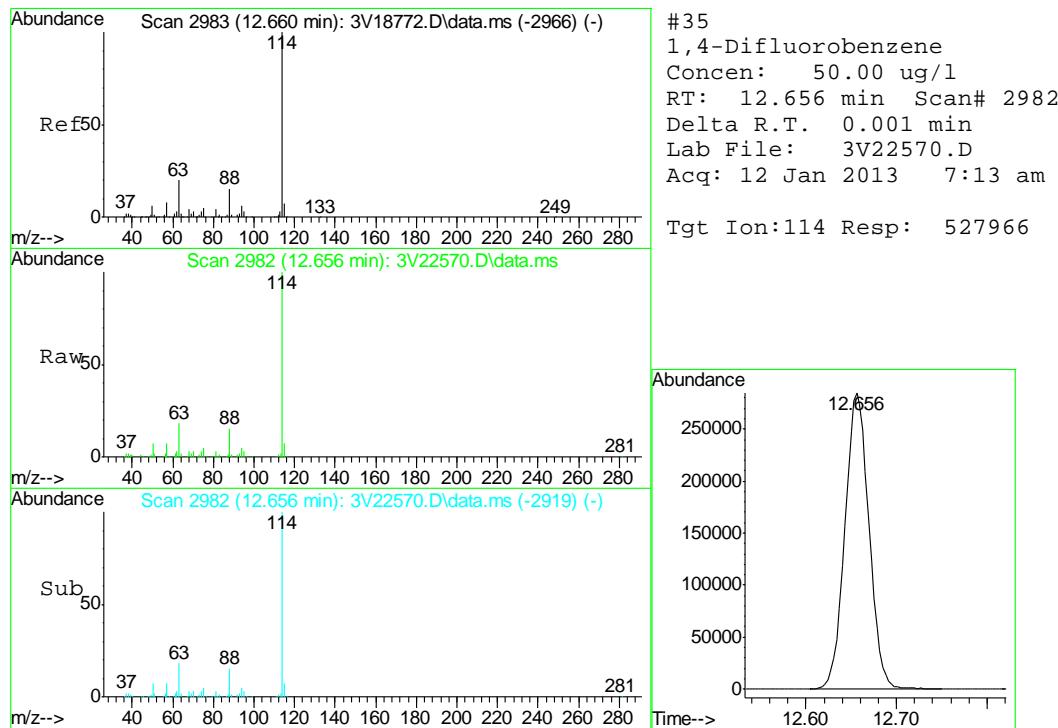
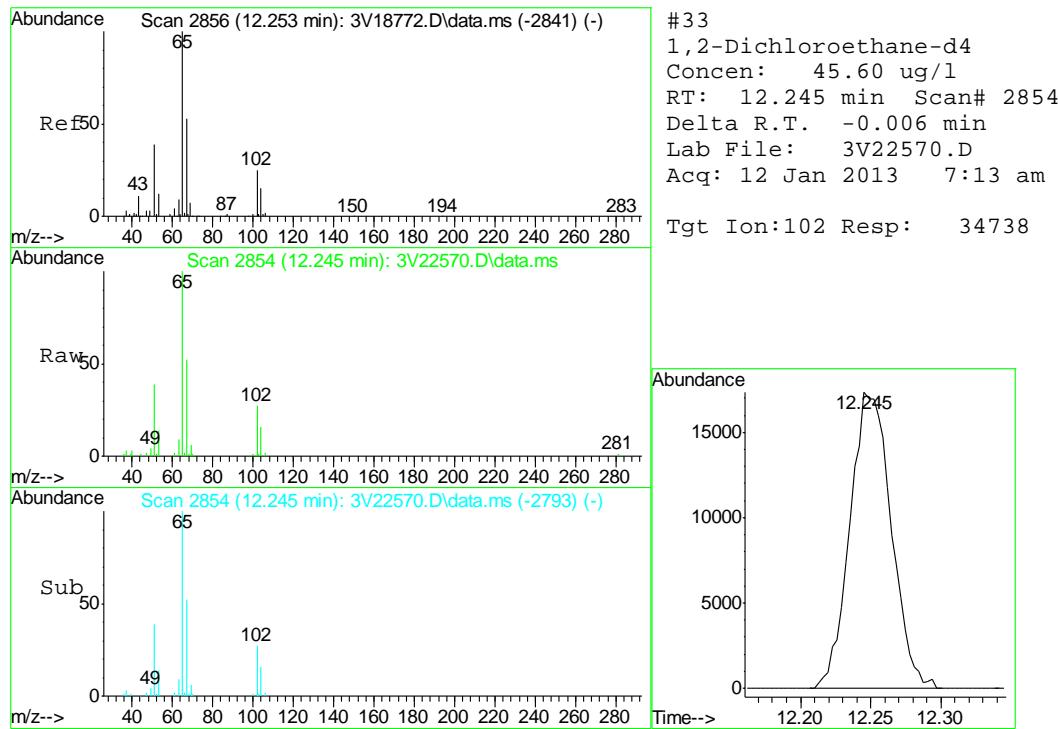
## Quantitation Report (QT Reviewed)

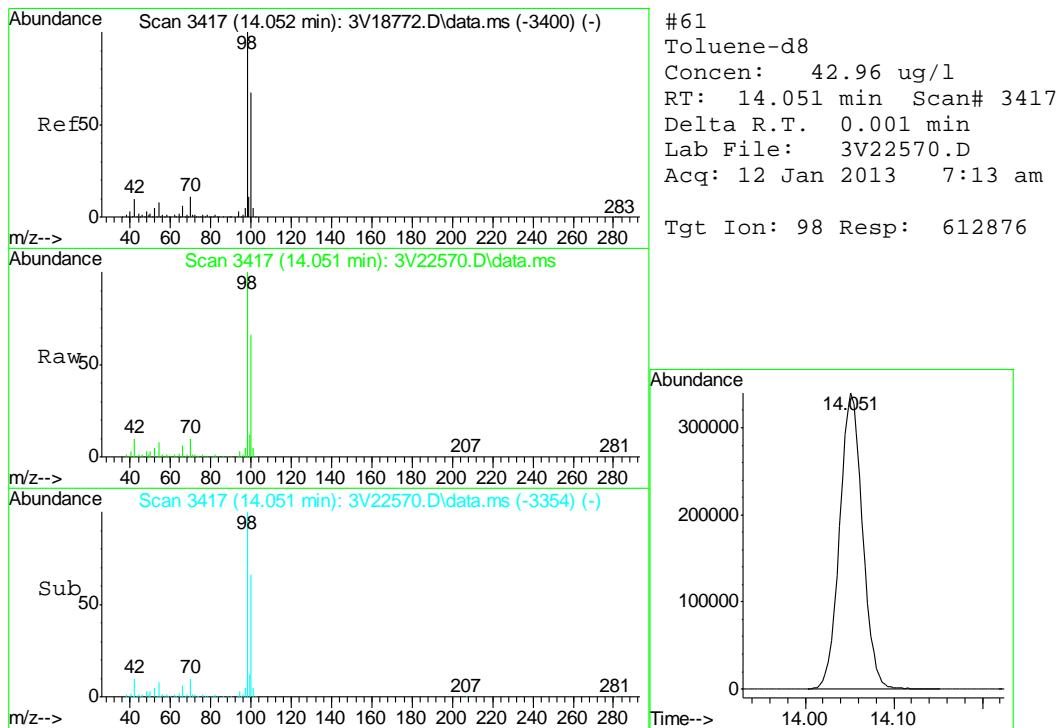
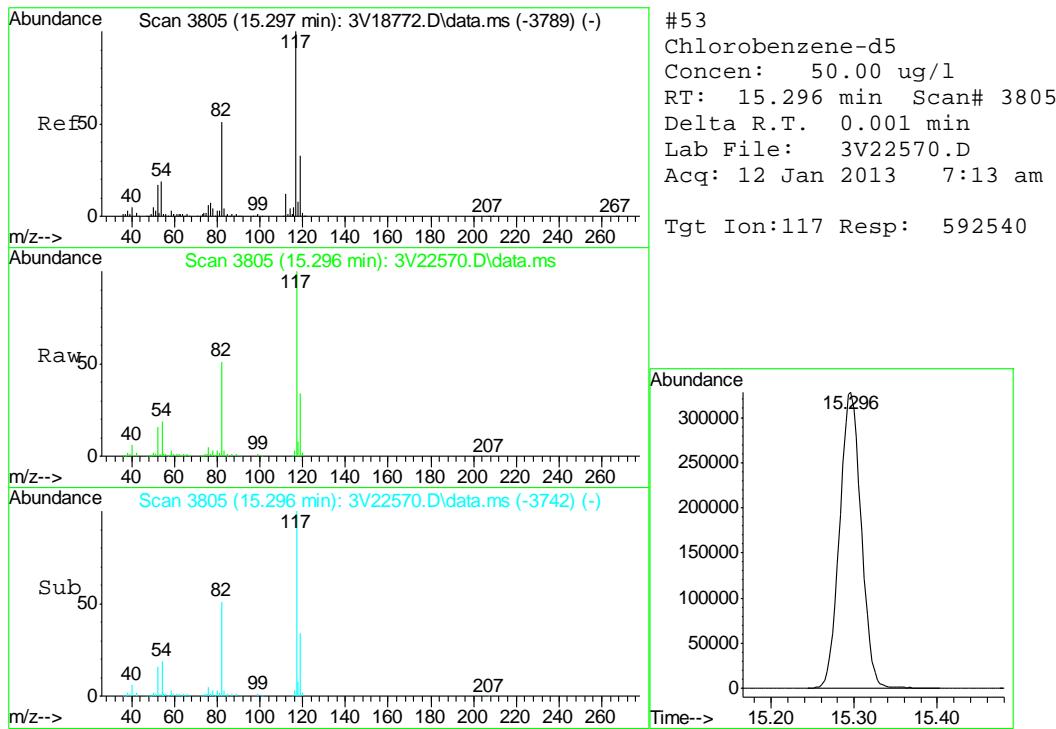
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 Data File : 3V22570.D  
 Acq On : 12 Jan 2013 7:13 am  
 Operator : BRETD  
 Sample : D42556-1  
 Misc : MS5218,V3V1327,5.066,,100,5,1  
 ALS Vial : 37 Sample Multiplier: 1

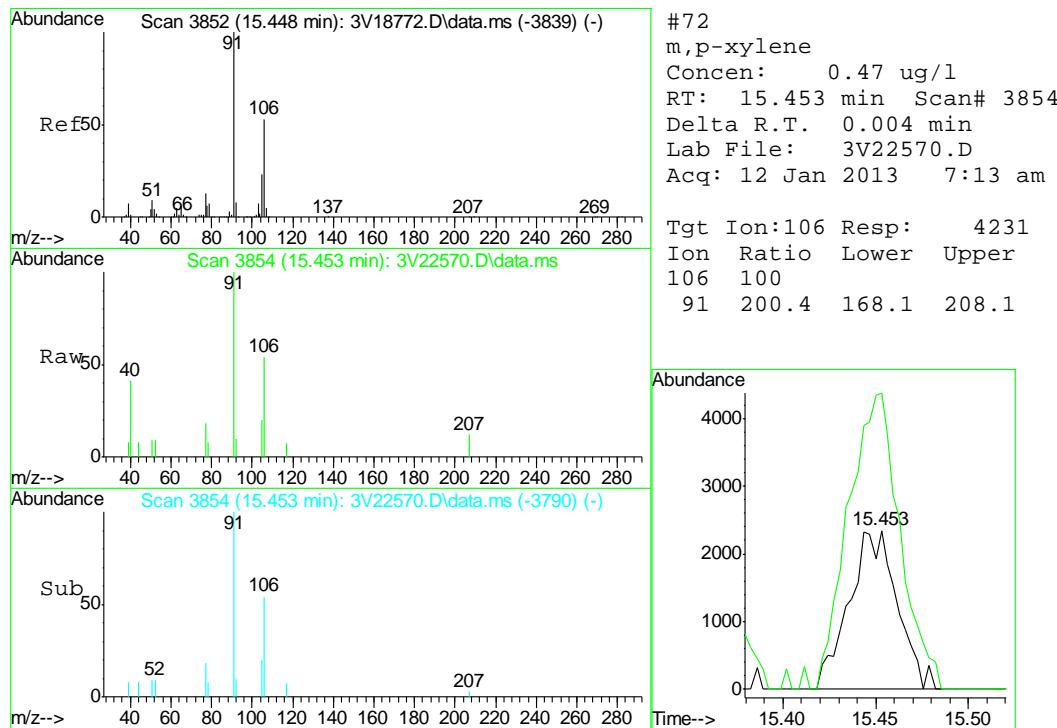
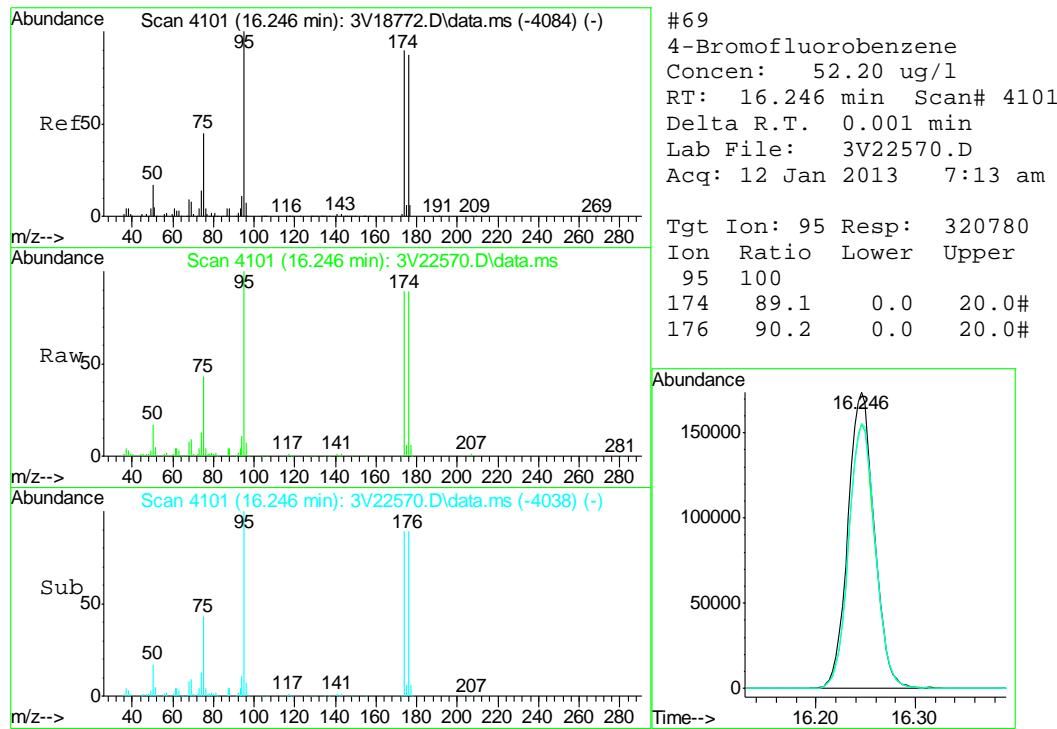
Quant Time: Jan 15 08:28:51 2013  
 Quant Method : C:\msdchem\1\METHODS\V3AP1299TVH1299SOIL.M  
 Quant Title : 8260  
 QLast Update : Thu Jan 03 11:40:16 2013  
 Response via : Initial Calibration

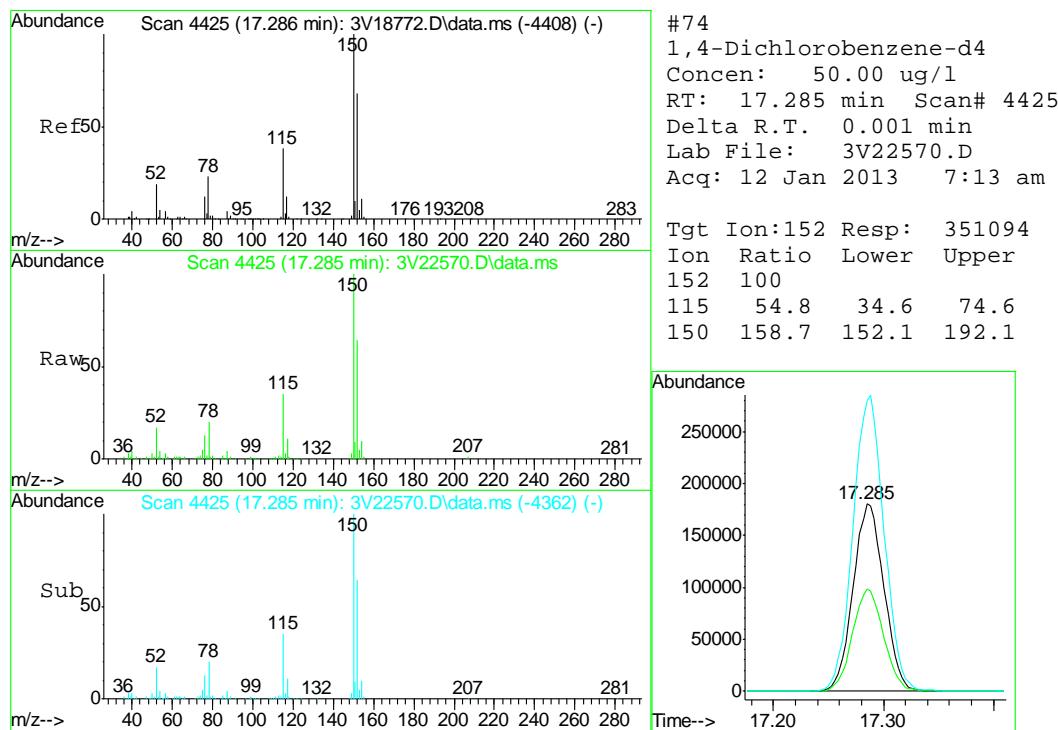
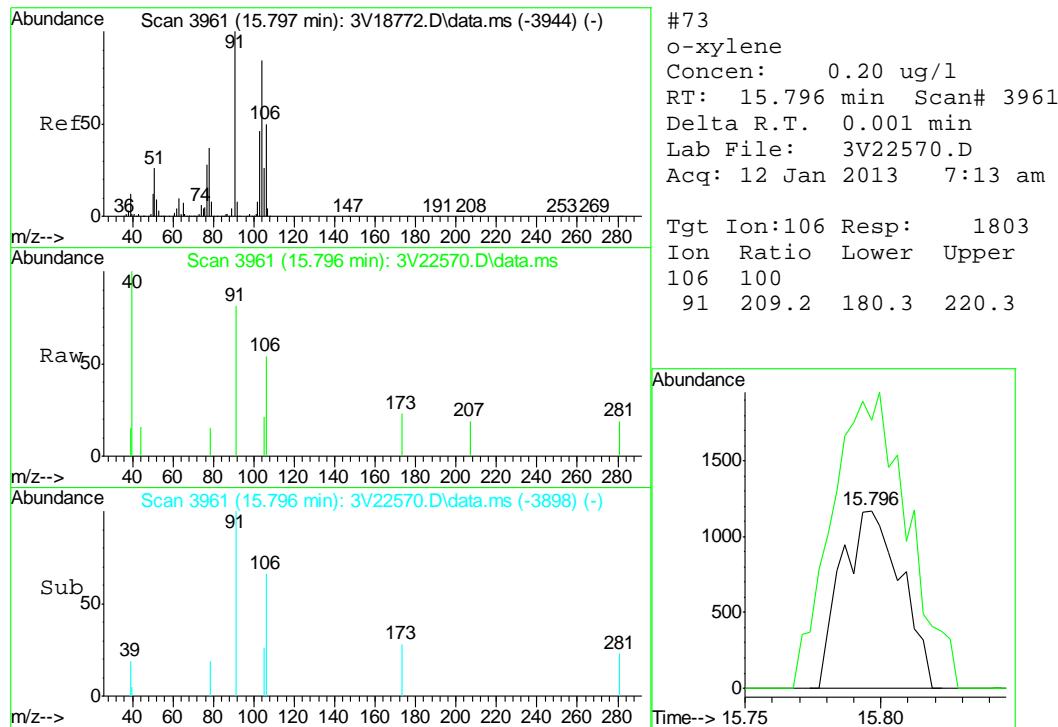


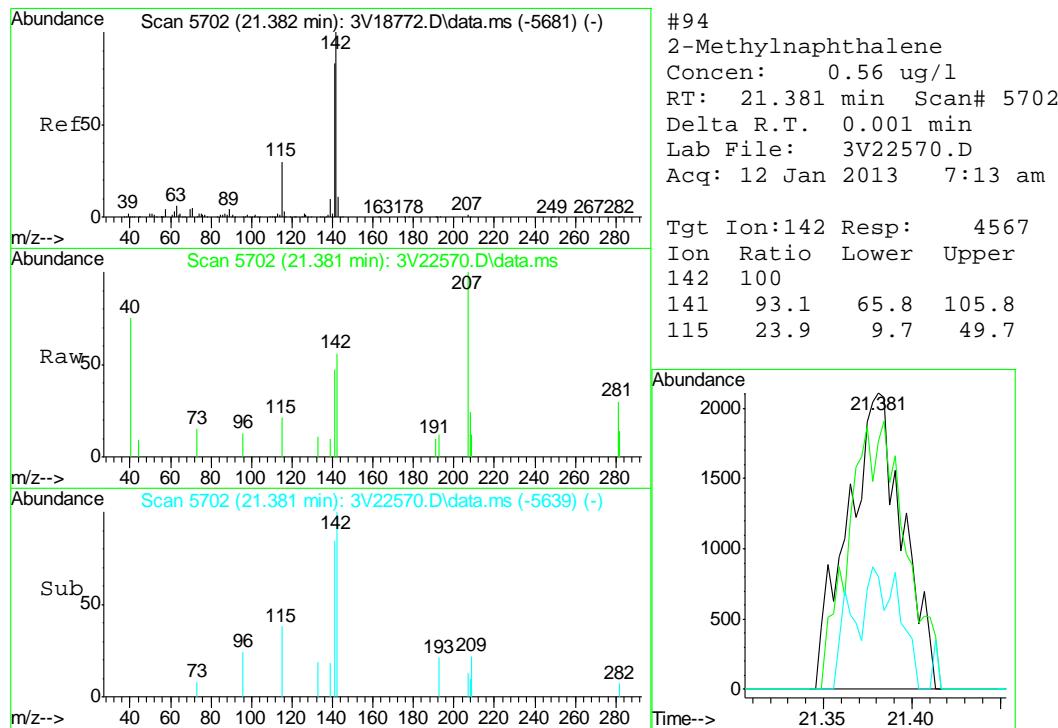
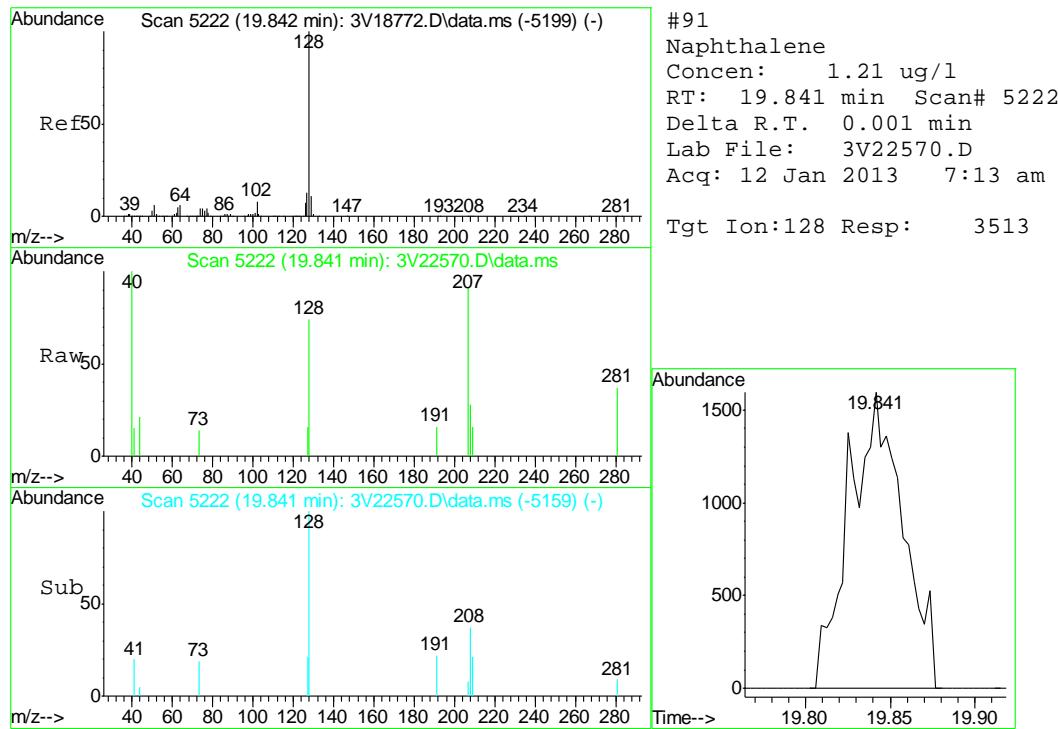


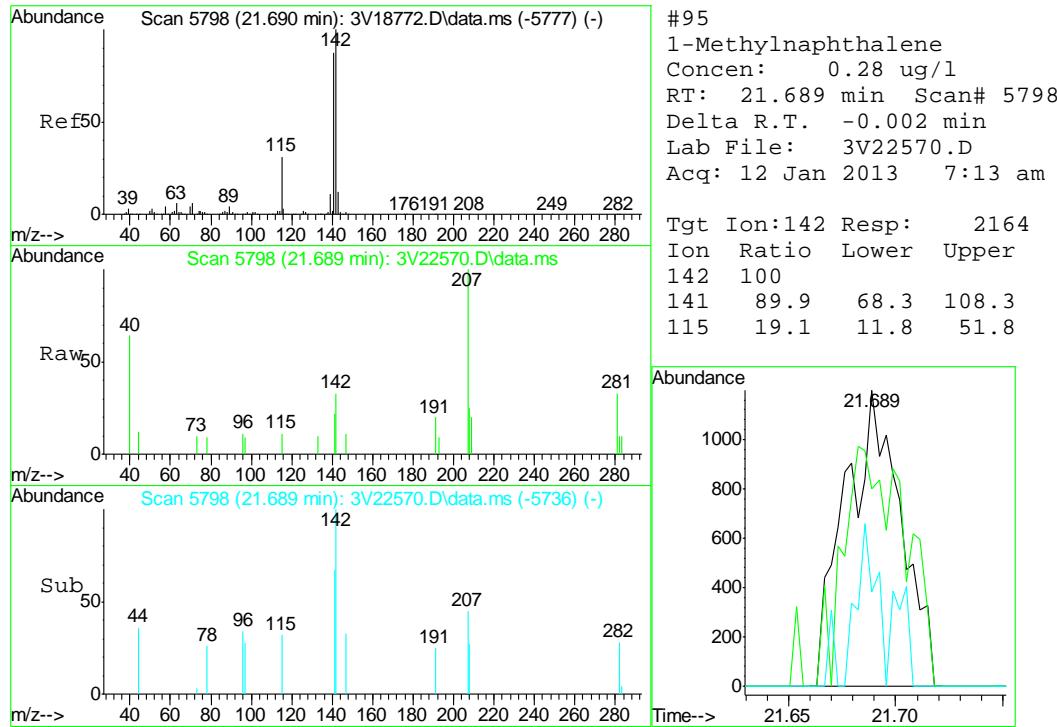












## Quantitation Report (QT Reviewed)

Manual Integrations  
APPROVED  
(compounds with "m" flag)

Judy Nelson  
01/15/13 10:36

Data Path : C:\msdchem\1\DATA\V3011113.S\  
Data File : 3V22571.D  
Acq On : 12 Jan 2013 7:45 am  
Operator : BRETD  
Sample : D42556-2  
Misc : MS5218,V3V1327,5.060,,100,5,1  
ALS Vial : 38 Sample Multiplier: 1

Quant Time: Jan 15 08:31:05 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	11.859	168	333289	50.00	ug/l	0.00
35) 1,4-Difluorobenzene	12.655	114	530134	50.00	ug/l	0.00
53) Chlorobenzene-d5	15.295	117	595022	50.00	ug/l	0.00
74) 1,4-Dichlorobenzene-d4	17.283	152	357062	50.00	ug/l	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
33) 1,2-Dichloroethane-d4	12.250	102	32371	43.40	ug/l	0.00
Spiked Amount 50.000	Range 70 - 130		Recovery	=	86.80%	
61) Toluene-d8	14.050	98	620064	43.28	ug/l	0.00
Spiked Amount 50.000	Range 70 - 130		Recovery	=	86.56%	
69) 4-Bromofluorobenzene	16.244	95	324673	52.61	ug/l	0.00
Spiked Amount 50.000	Range 70 - 130		Recovery	=	105.22%	

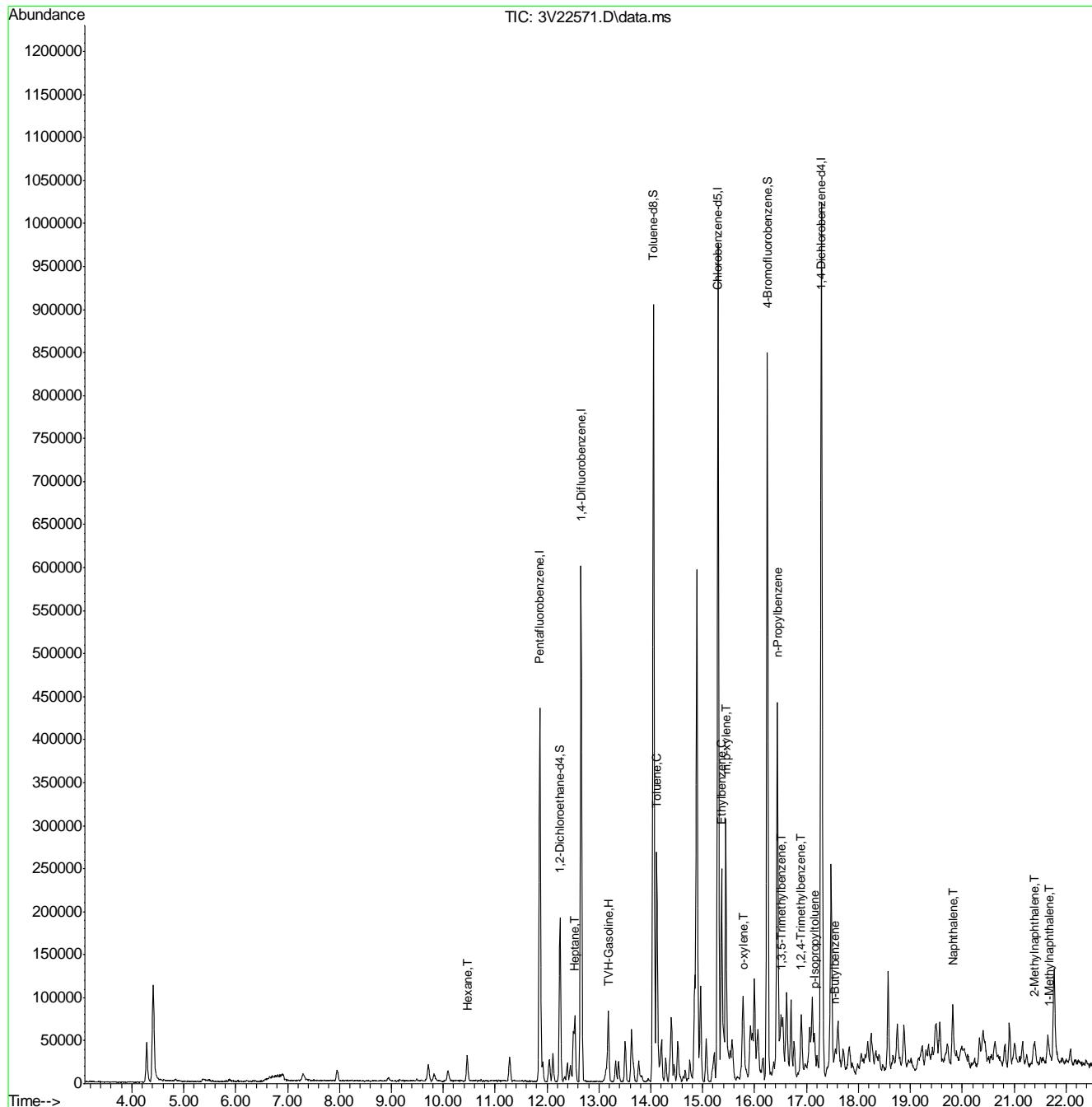
Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
1) TVH-Gasoline	13.200	TIC	7998907m	272.57	ug/l	
41) Hexane	10.460	57	14962	2.41	ug/l	100
43) Heptane	12.536	43	30804	4.33	ug/l	92
62) Toluene	14.114	92	68281	5.53	ug/l	97
66) Ethylbenzene	15.362	91	54248	2.50	ug/l	97
72) m,p-xylene	15.445	106	97828	10.71	ug/l	99
73) o-xylene	15.798	106	10886	1.23	ug/l	99
77) n-Propylbenzene	16.427	91	15879	0.65	ug/l	100
80) 1,3,5-Trimethylbenzene	16.510	105	29541m	1.61	ug/l	
82) 1,2,4-Trimethylbenzene	16.898	105	24027	1.28	ug/l	91
86) p-Isopropyltoluene	17.155	119	23185	1.11	ug/l	96
88) n-Butylbenzene	17.540	91	8553	0.50	ug/l	# 73
91) Naphthalene	19.833	128	10434	1.57	ug/l	100
94) 2-Methylnaphthalene	21.376	142	15889	1.92	ug/l	94
95) 1-Methylnaphthalene	21.691	142	6363	0.81	ug/l	96

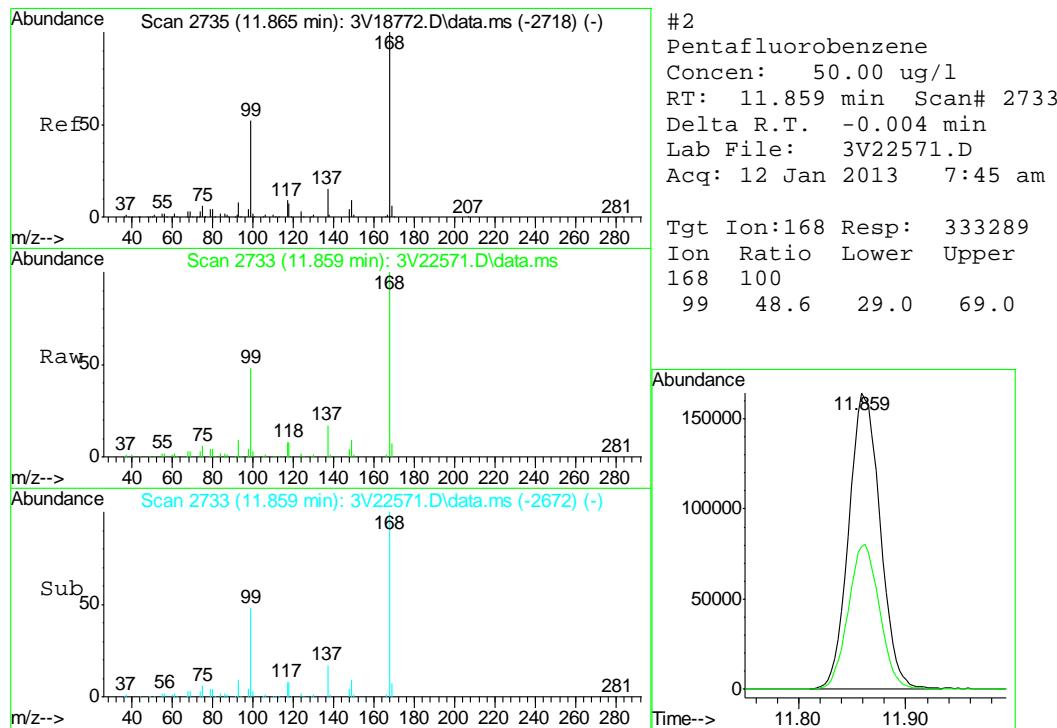
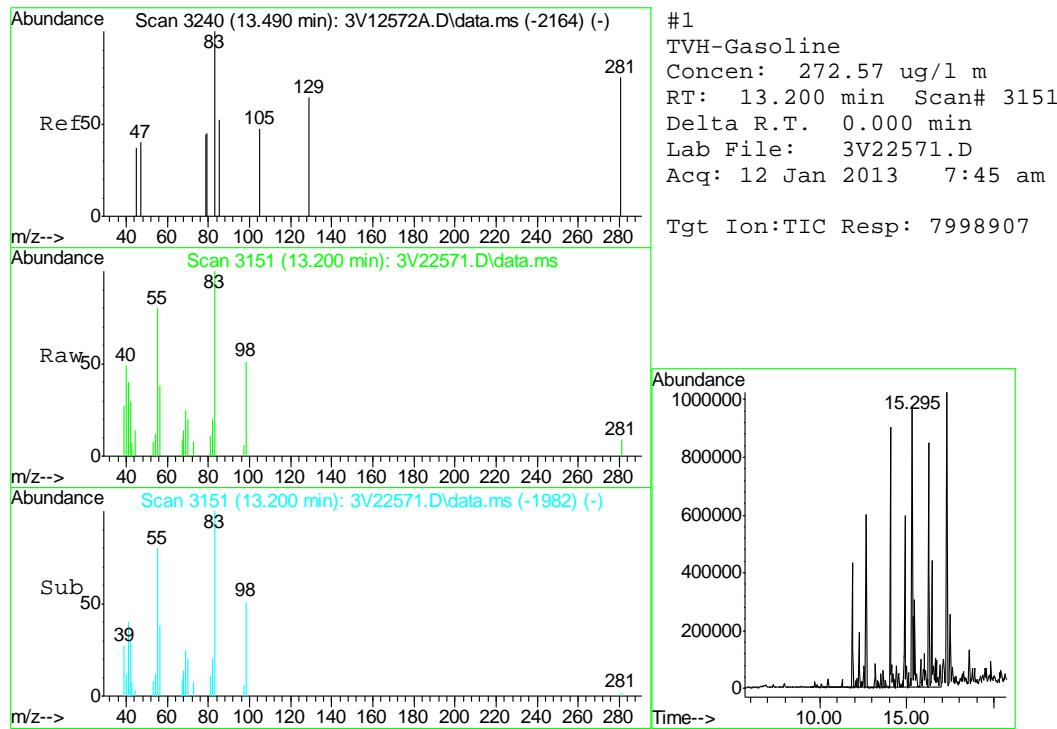
(#) = qualifier out of range (m) = manual integration (+) = signals summed

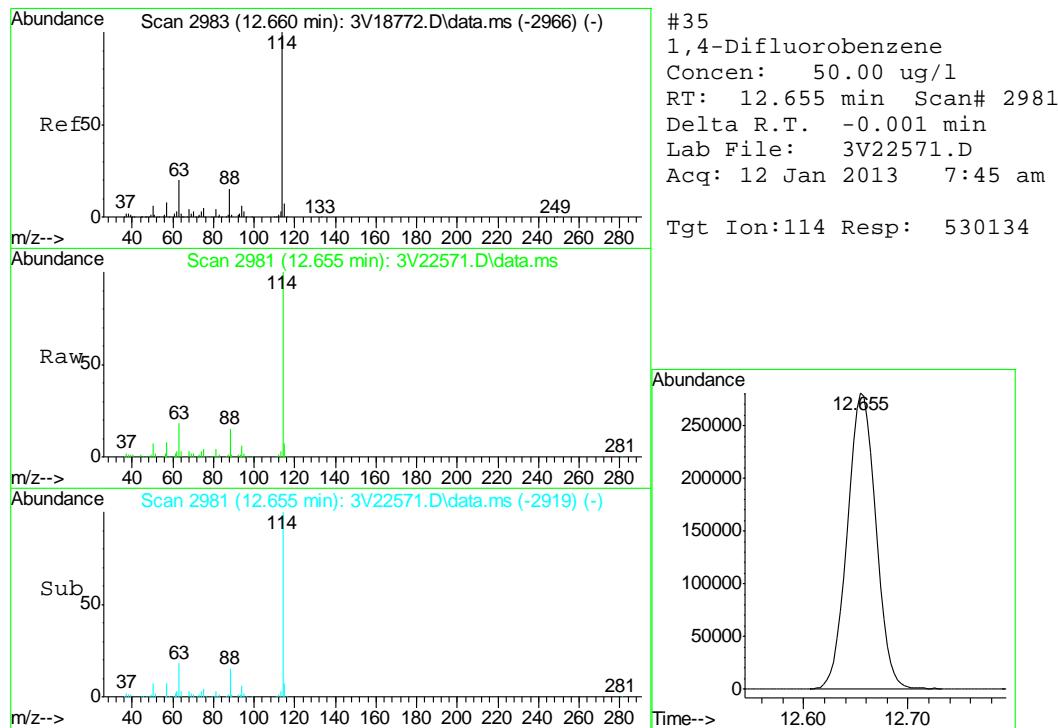
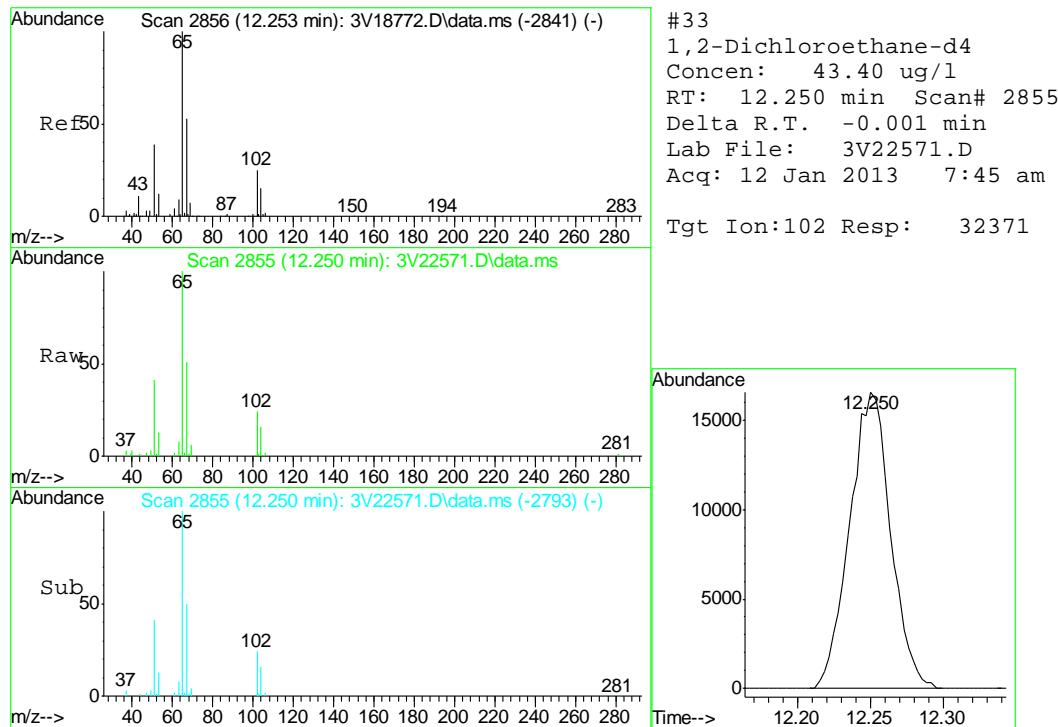
## Quantitation Report (QT Reviewed)

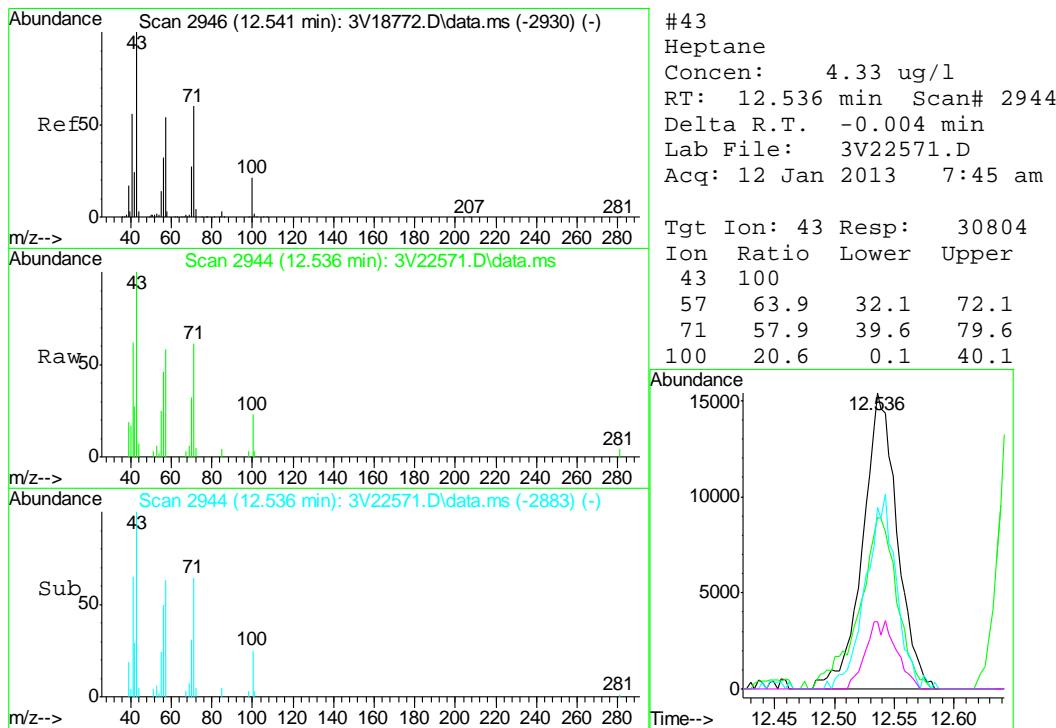
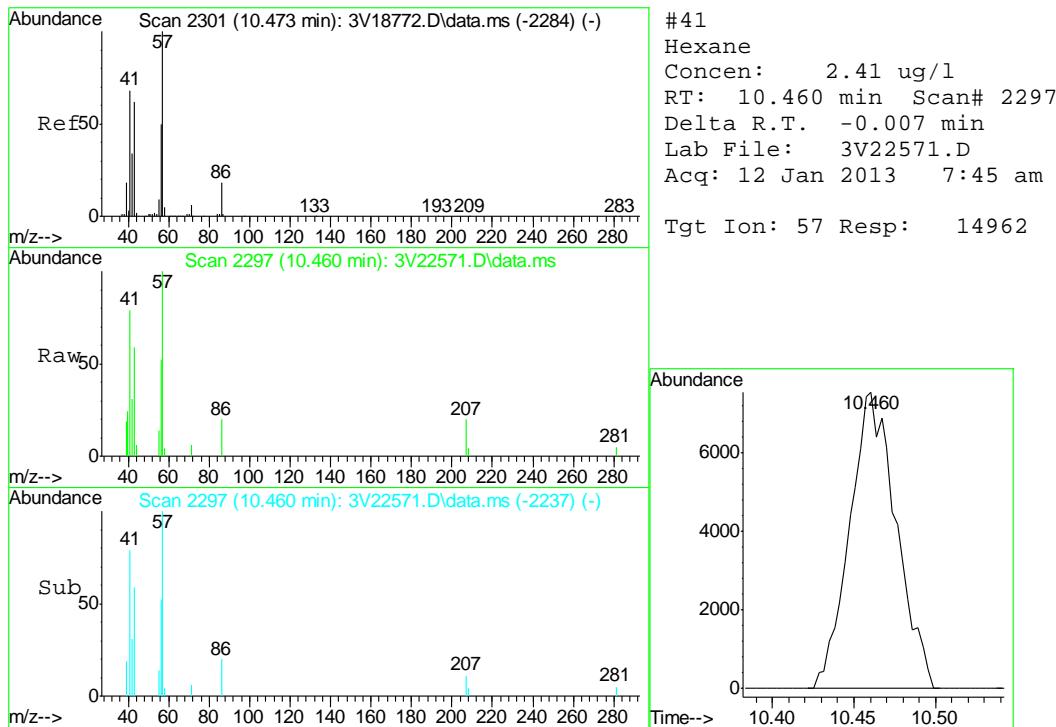
Data Path : C:\msdchem\1\DATA\V3011113.S\  
 Data File : 3V22571.D  
 Acq On : 12 Jan 2013 7:45 am  
 Operator : BRETD  
 Sample : D42556-2  
 Misc : MS5218,V3V1327,5.060,,100,5,1  
 ALS Vial : 38 Sample Multiplier: 1

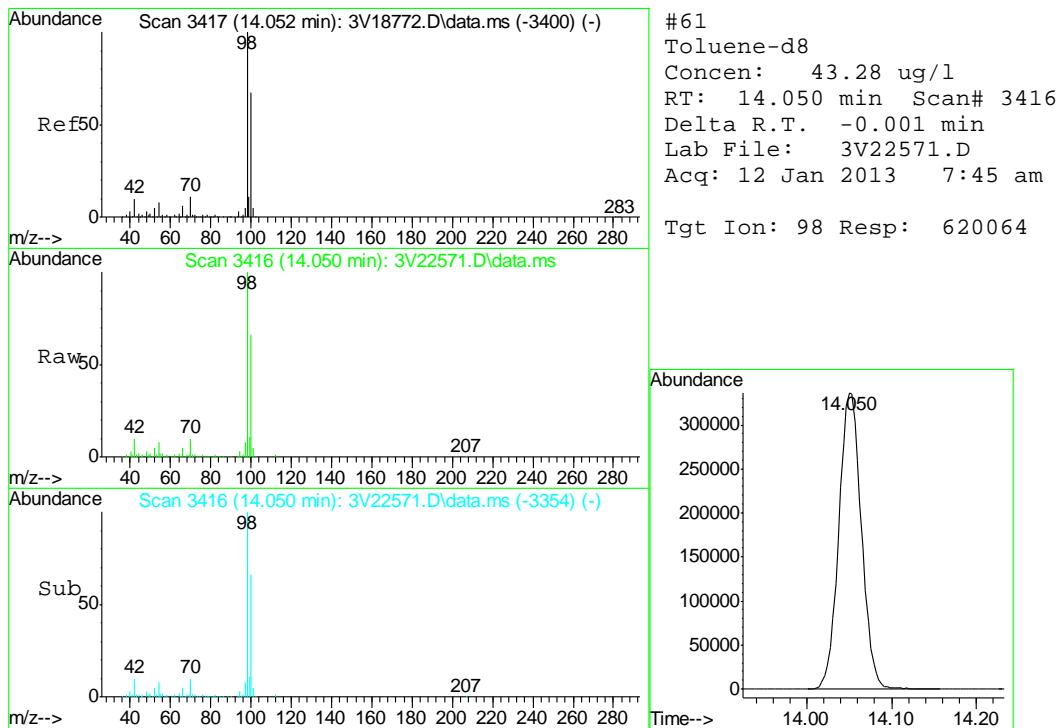
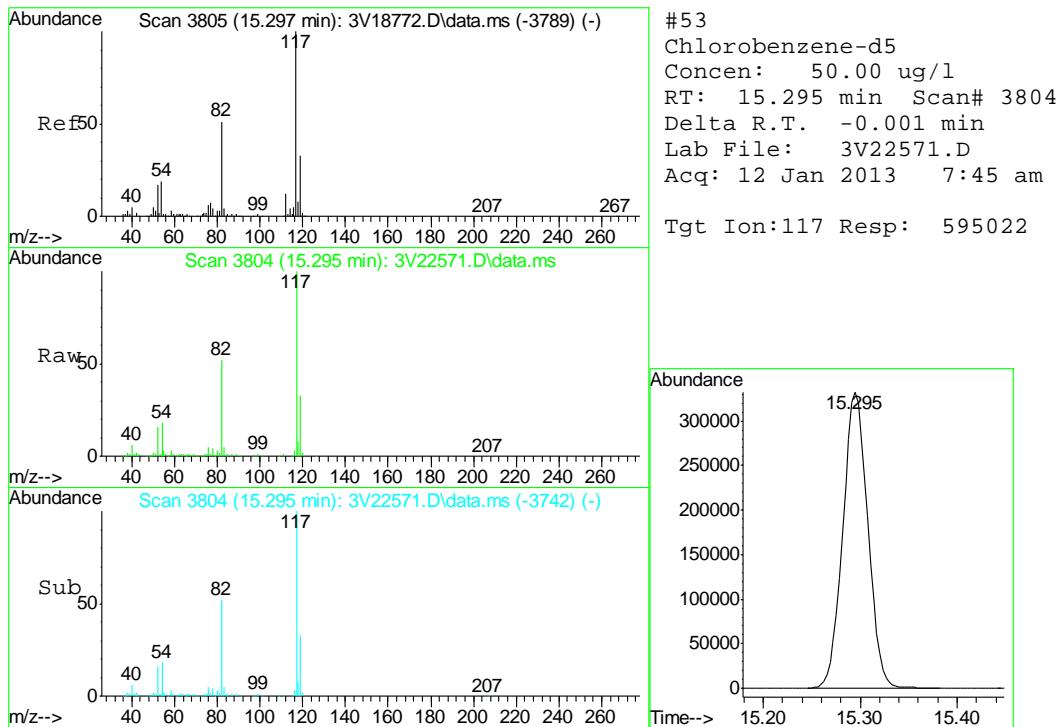
Quant Time: Jan 15 08:31:05 2013  
 Quant Method : C:\msdchem\1\METHODS\V3AP1299TVH1299SOIL.M  
 Quant Title : 8260  
 QLast Update : Thu Jan 03 11:40:16 2013  
 Response via : Initial Calibration

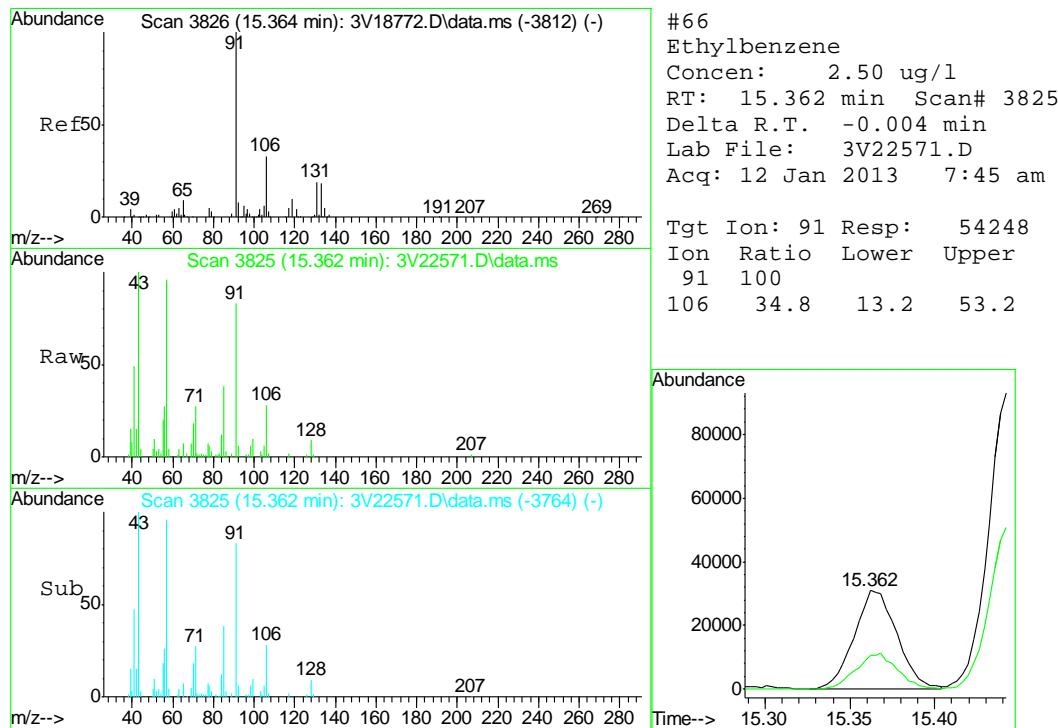
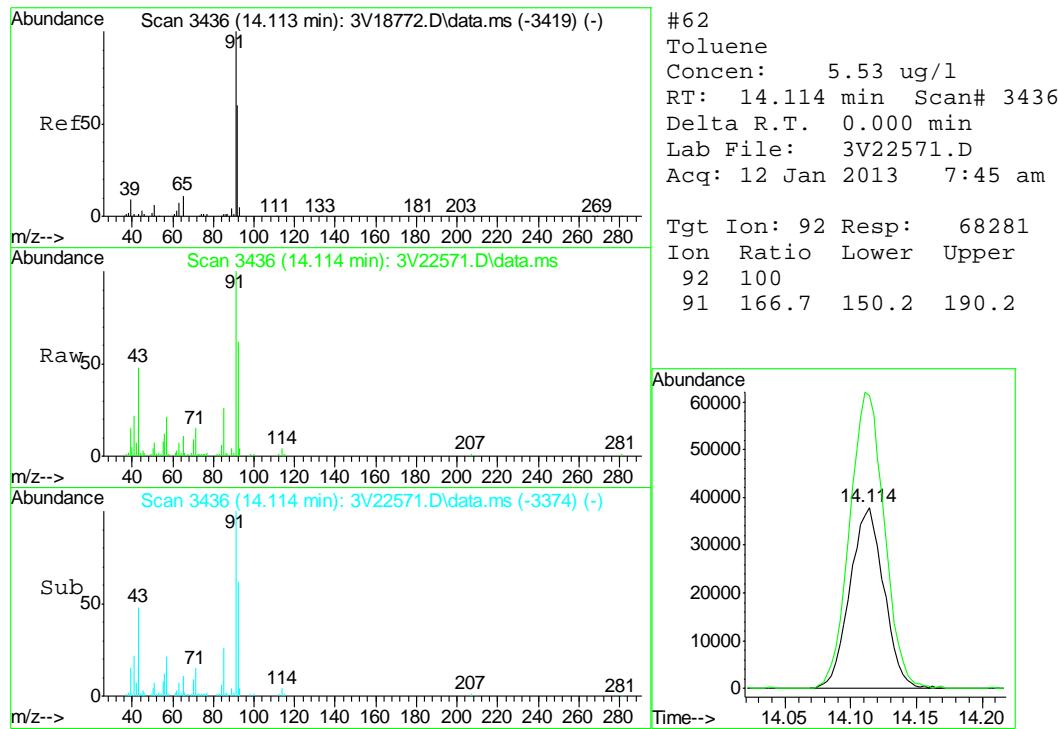


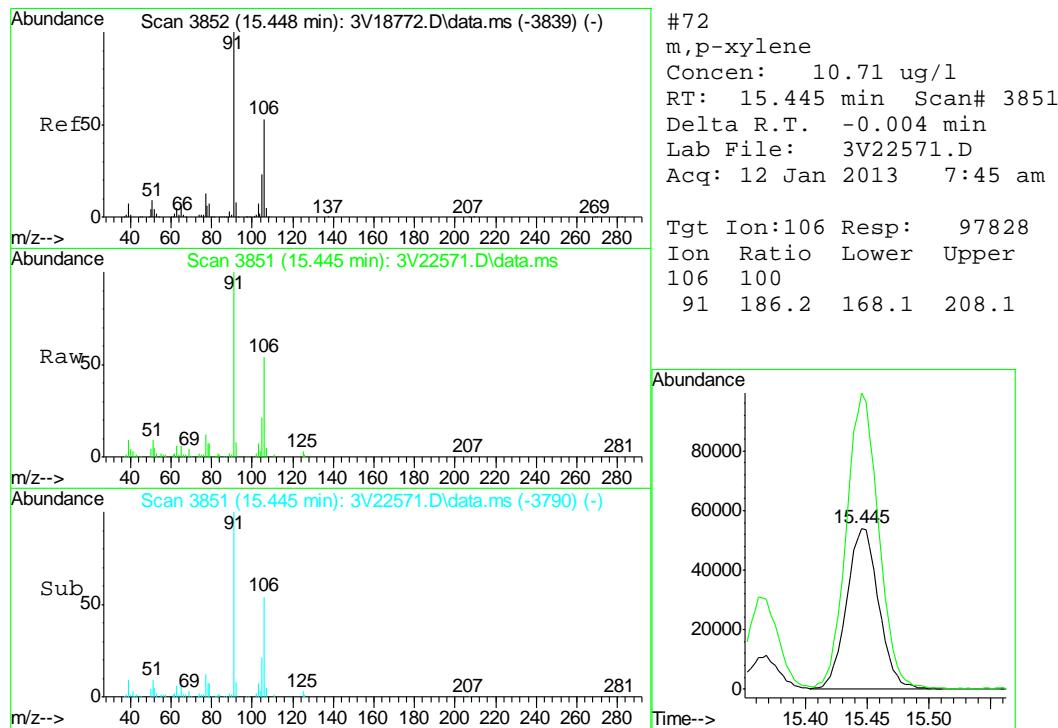
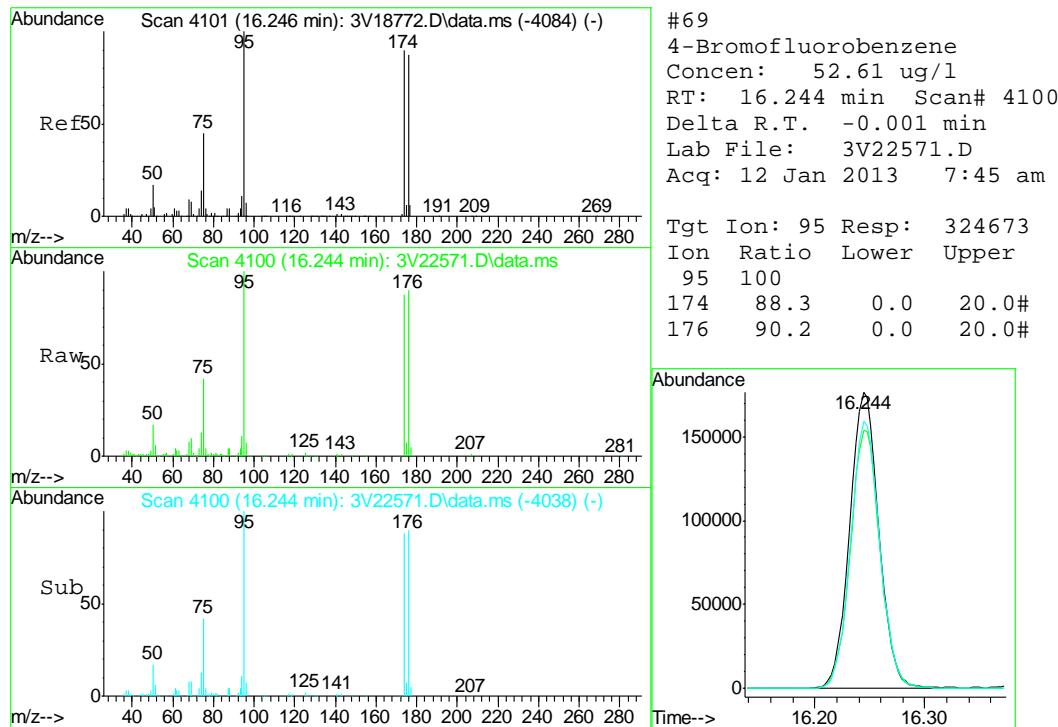


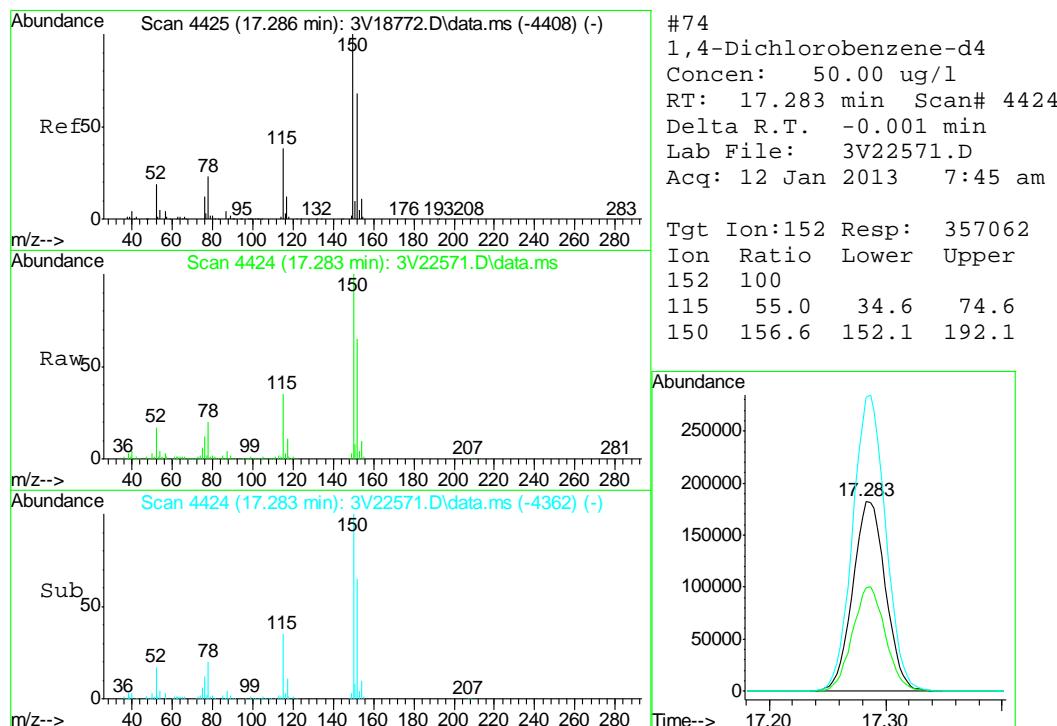
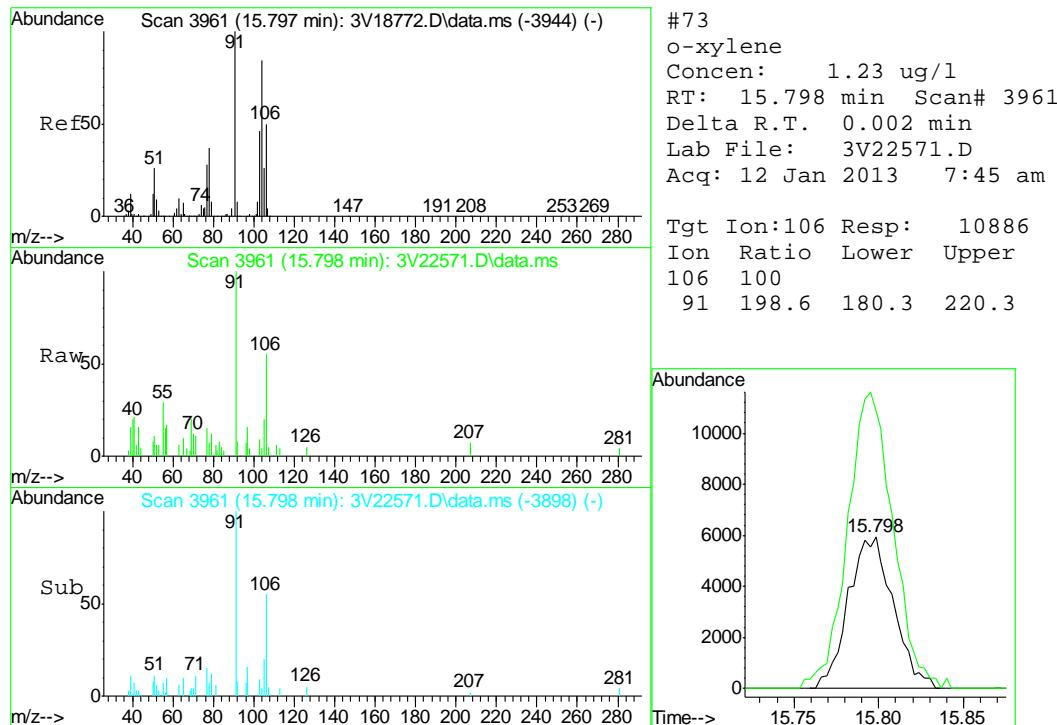


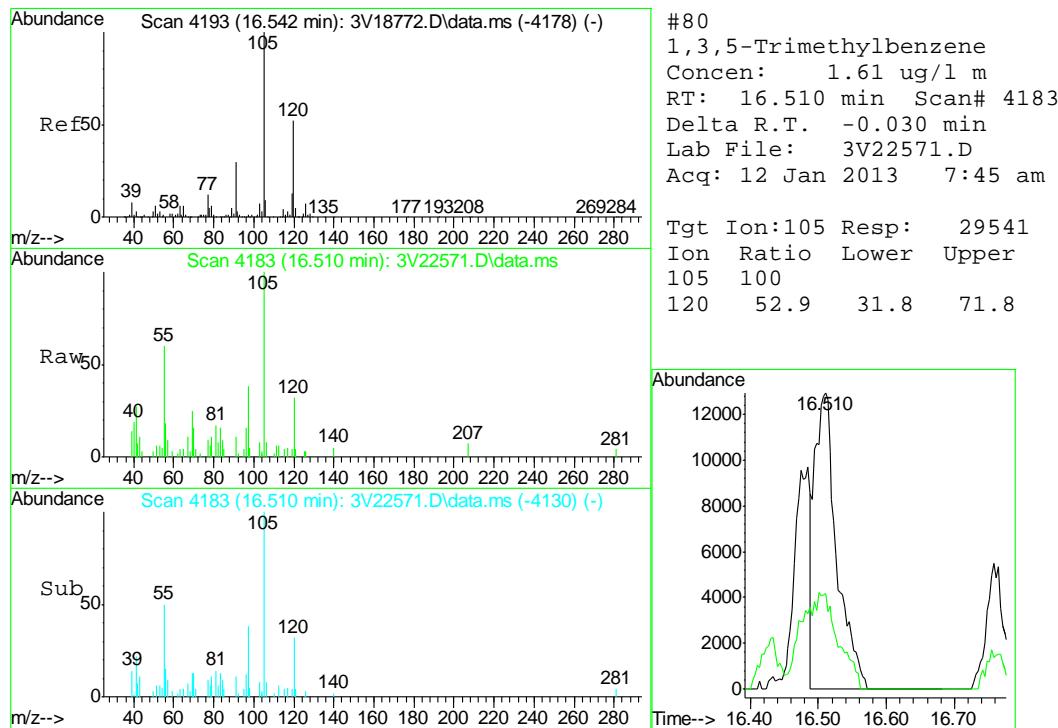
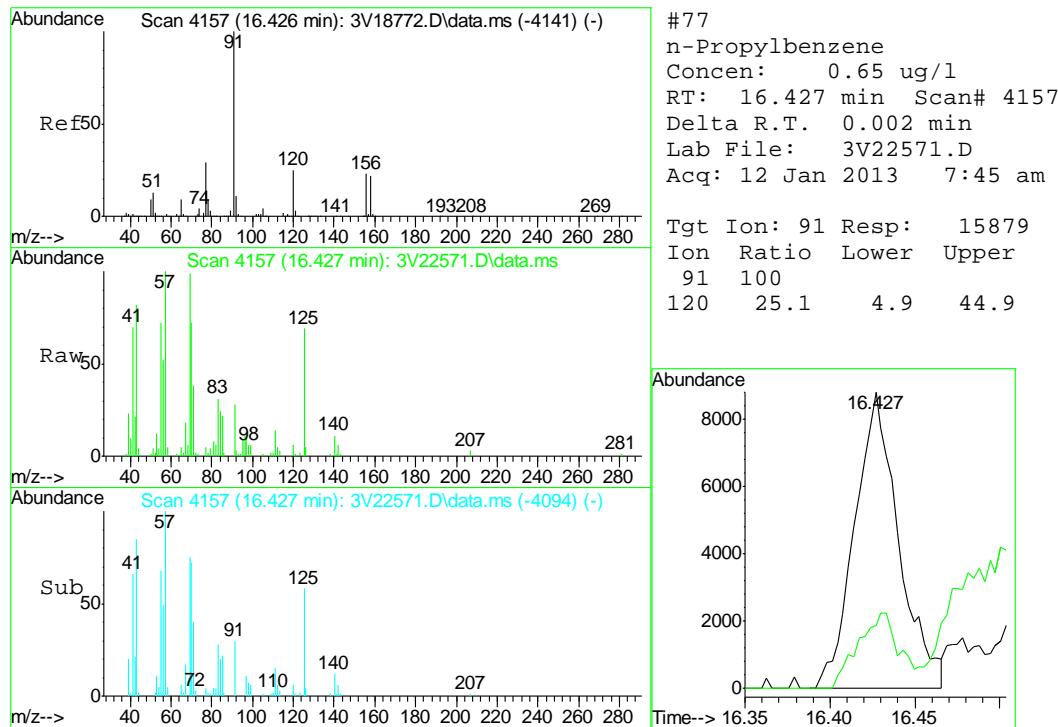


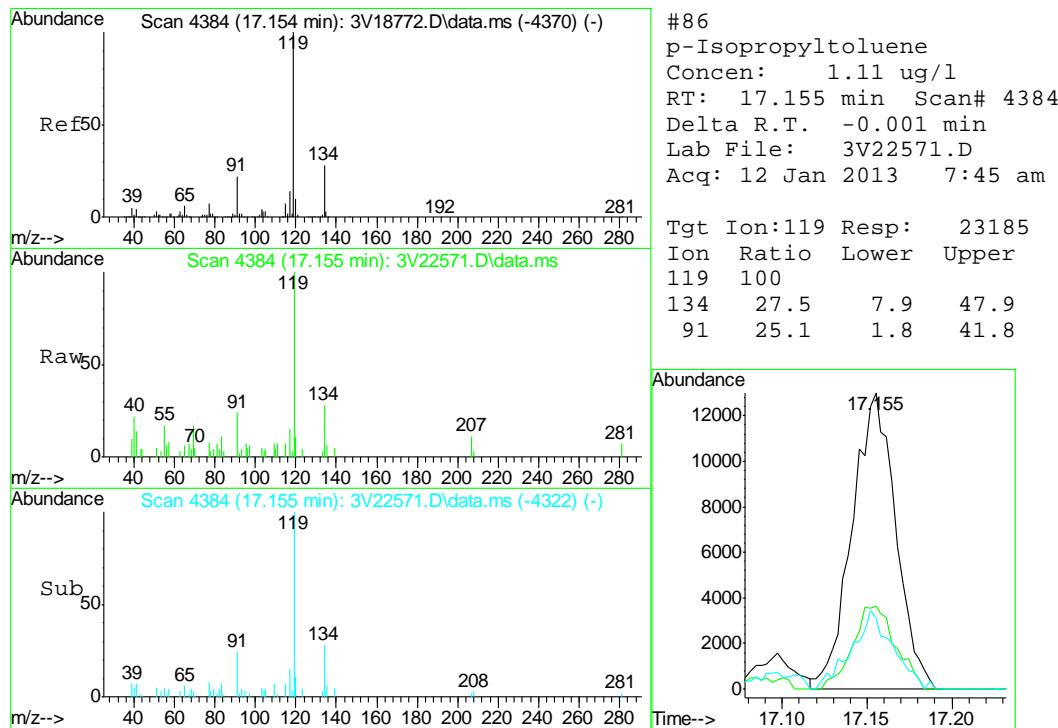
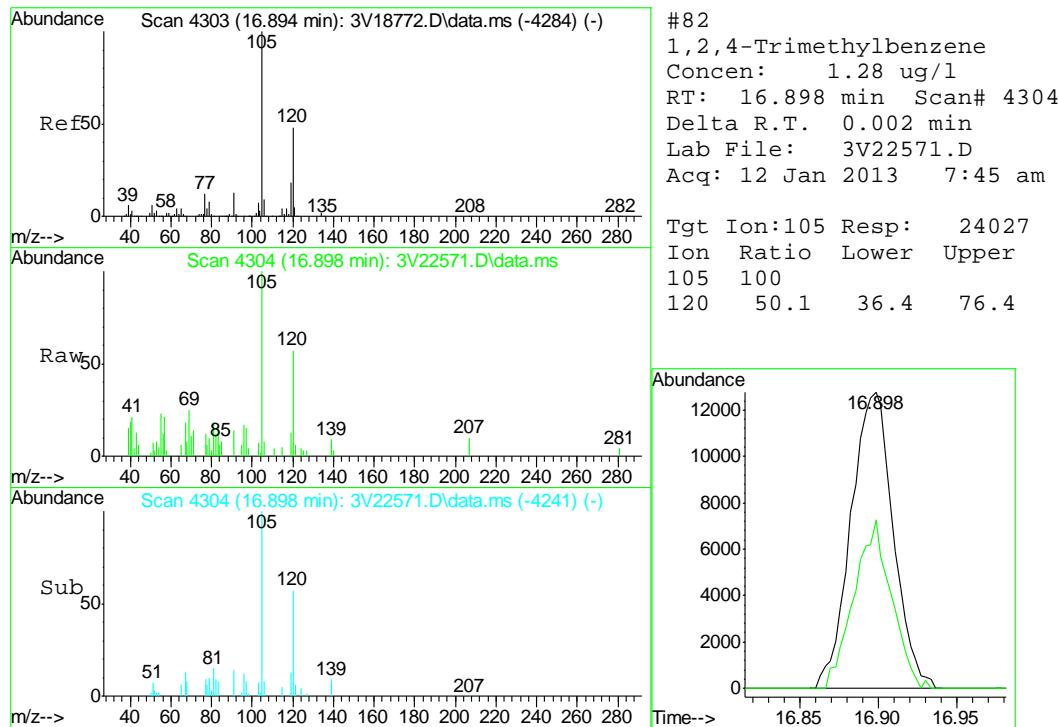


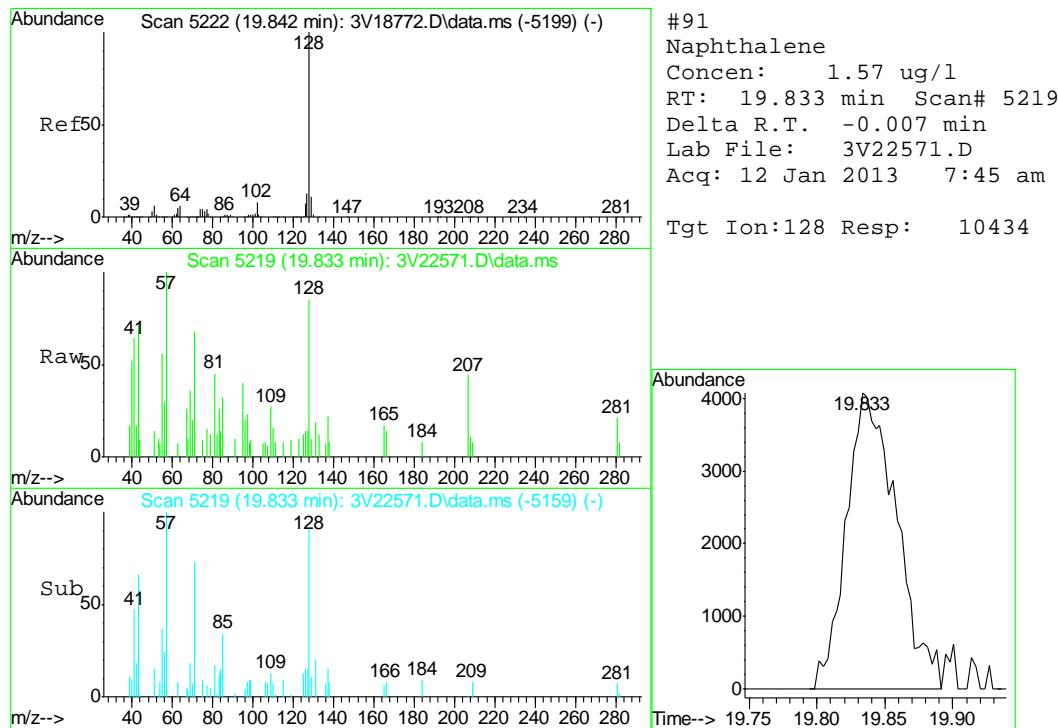
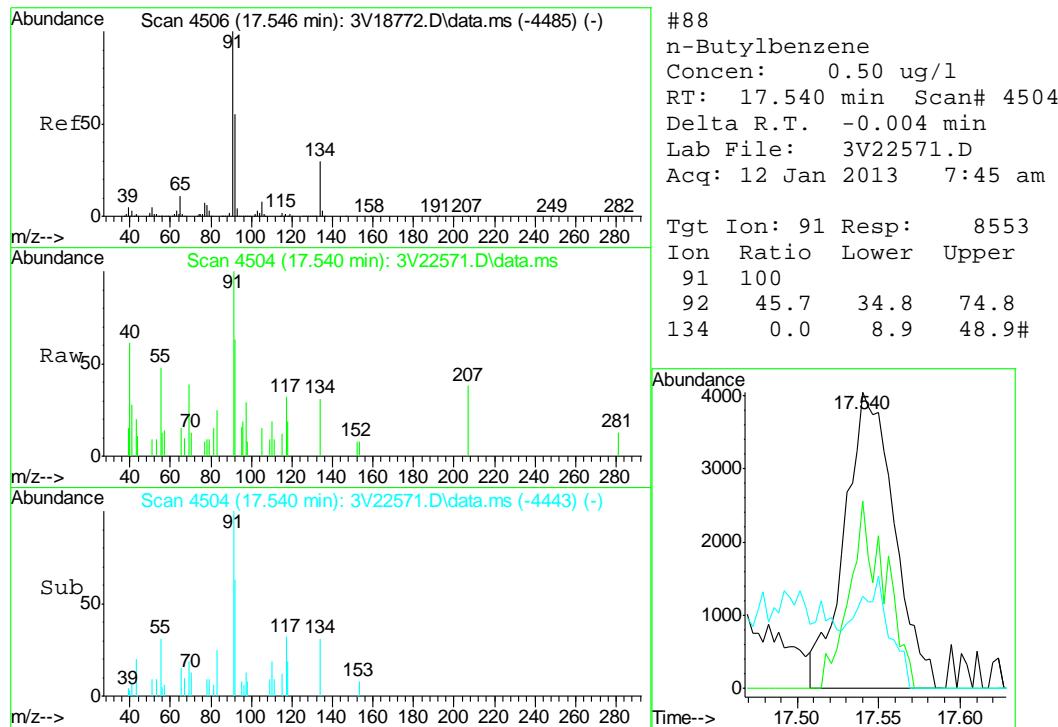


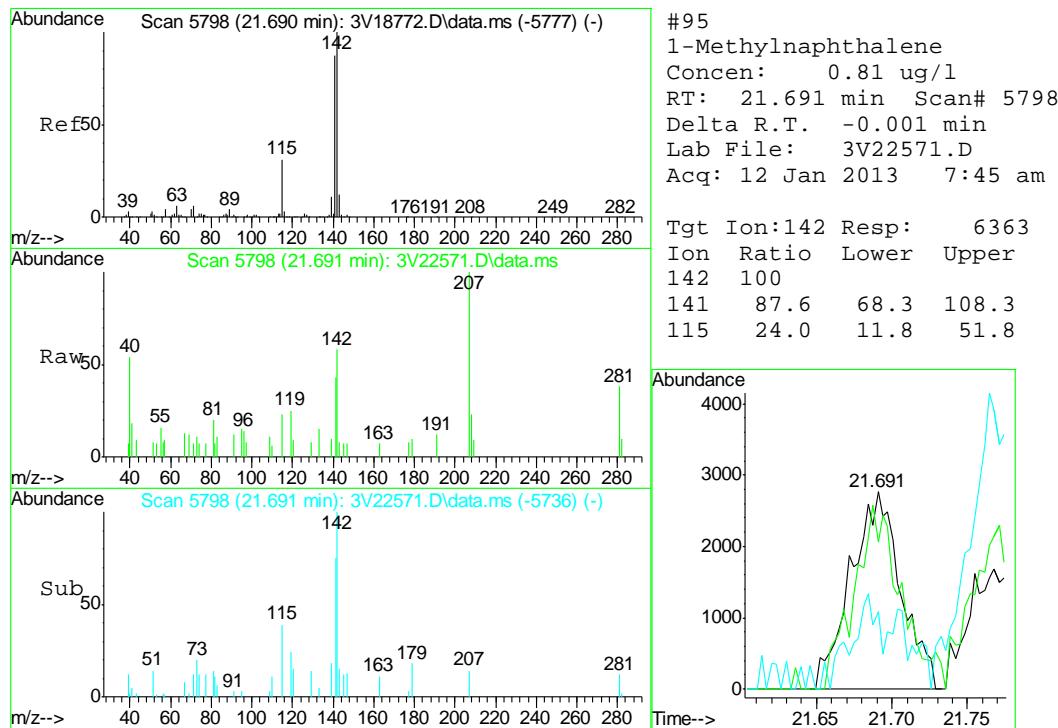
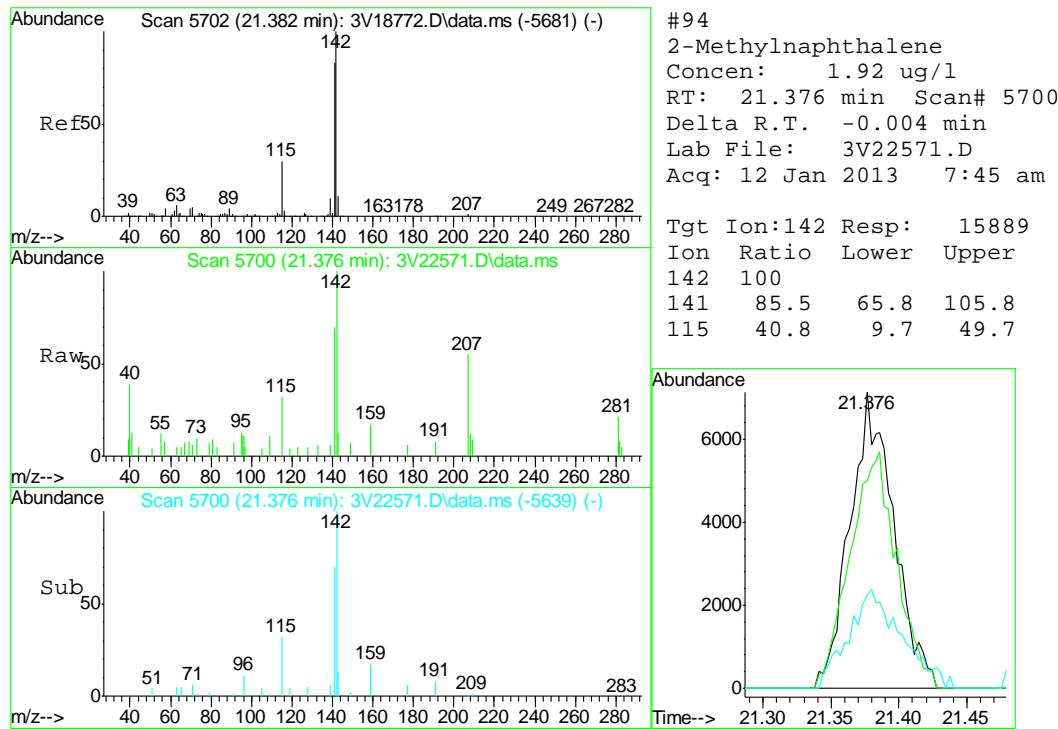












## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\V3011113.S\  
 Data File : 3V22560.D  
 Acq On : 12 Jan 2013 2:00 am  
 Operator : BRETD  
 Sample : MB  
 Misc : MS5218,V3V1327,5.00,,100,5,1  
 ALS Vial : 27 Sample Multiplier: 1

Quant Time: Jan 15 08:11:29 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	308471	50.00	ug/l	0.00
35) 1,4-Difluorobenzene	12.659	114	500082	50.00	ug/l	0.00
53) Chlorobenzene-d5	15.295	117	525362	50.00	ug/l	0.00
74) 1,4-Dichlorobenzene-d4	17.287	152	304372	50.00	ug/l	0.00

System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	12.251	102	33686	48.79	ug/l	0.00
Spiked Amount 50.000	Range 70 - 130		Recovery	=	97.58%	
61) Toluene-d8	14.054	98	562082	44.44	ug/l	0.00
Spiked Amount 50.000	Range 70 - 130		Recovery	=	88.88%	
69) 4-Bromofluorobenzene	16.245	95	265181	48.67	ug/l	0.00
Spiked Amount 50.000	Range 70 - 130		Recovery	=	97.34%	

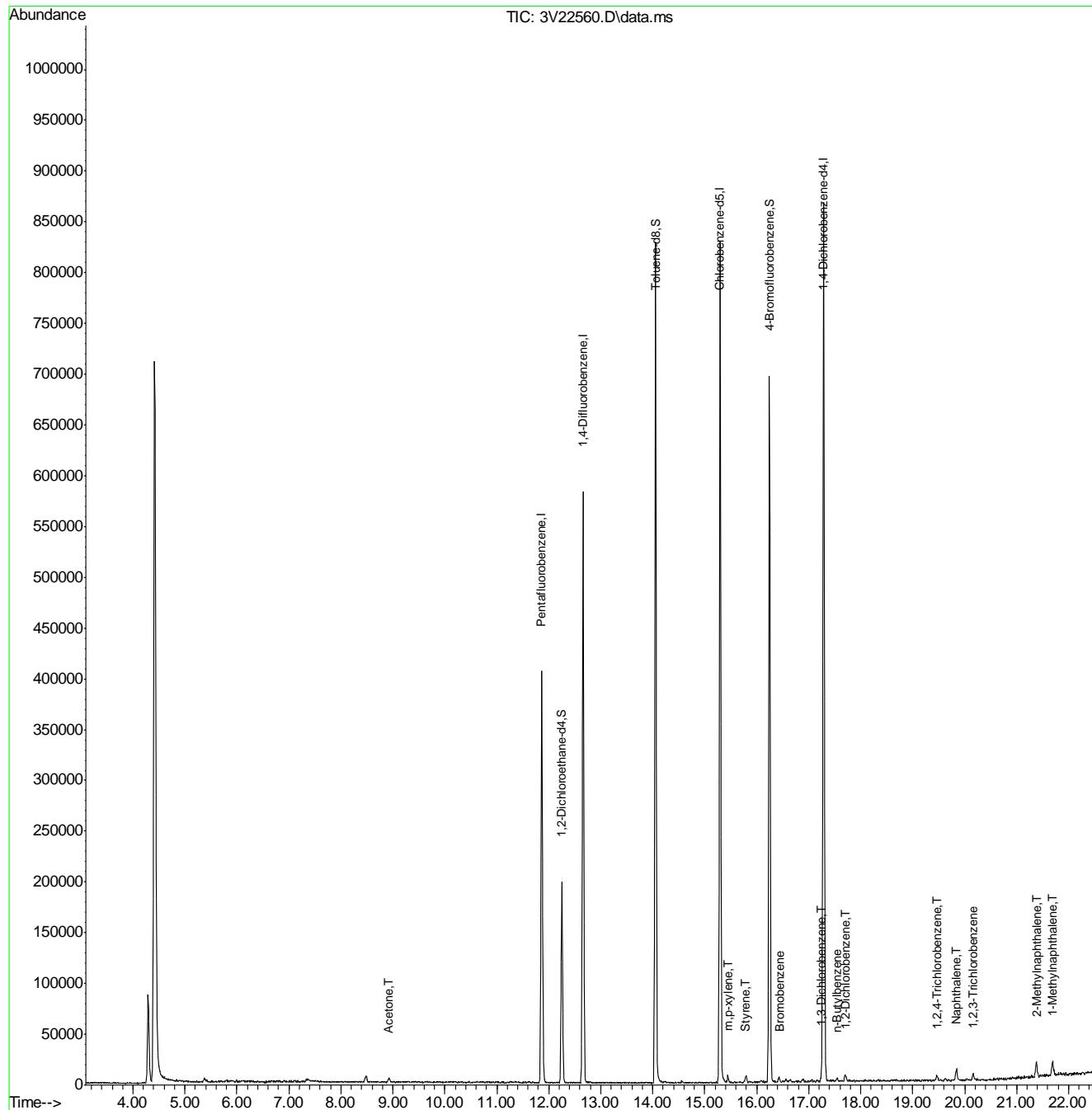
Target Compounds					Qvalue
15) Acetone	8.925	43	6973	0.73	ug/l 100
70) Bromobenzene	16.441	156	1174	0.21	ug/l 90
71) Styrene	15.796	104	1960	0.56	ug/l 86
72) m,p-xylene	15.453	106	2503	0.31	ug/l 90
84) 1,3-Dichlorobenzene	17.239	146	3019	0.28	ug/l 96
87) 1,2-Dichlorobenzene	17.704	146	3319	0.32	ug/l 96
88) n-Butylbenzene	17.541	91	3129	0.22	ug/l 88
90) 1,2,4-Trichlorobenzene	19.462	180	3118	0.48	ug/l 92
91) Naphthalene	19.838	128	16063	2.00	ug/l 100
93) 1,2,3-Trichlorobenzene	20.168	180	3642	0.57	ug/l 94
94) 2-Methylnaphthalene	21.384	142	13326	1.89	ug/l 97
95) 1-Methylnaphthalene	21.692	142	11465	1.71	ug/l 98

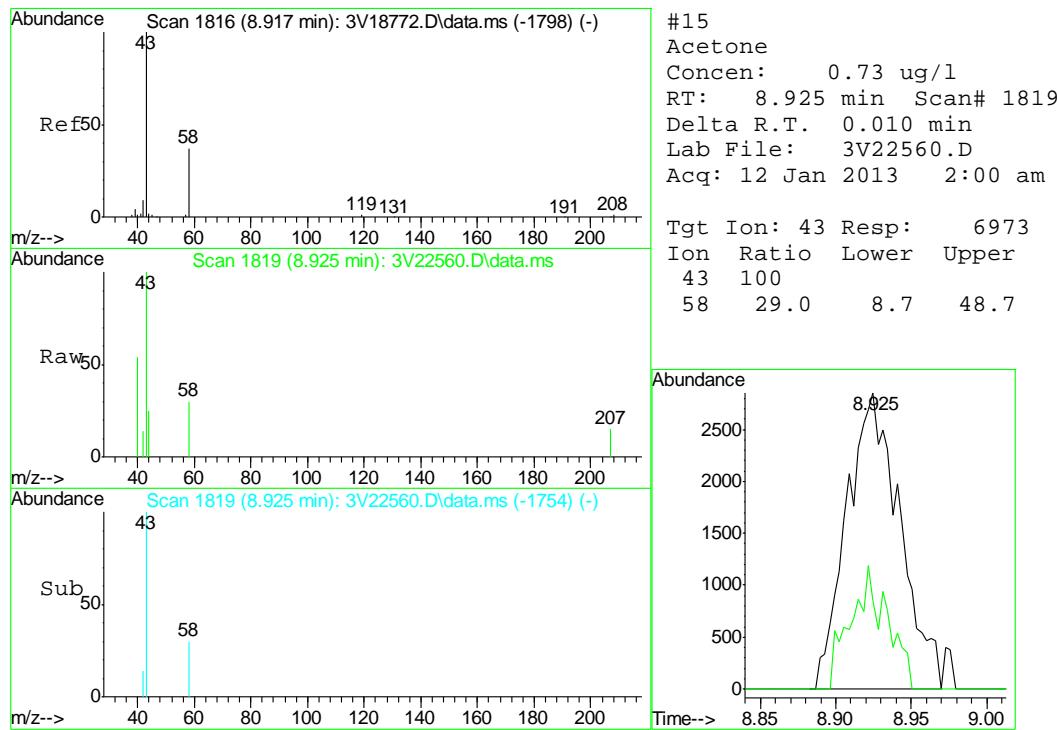
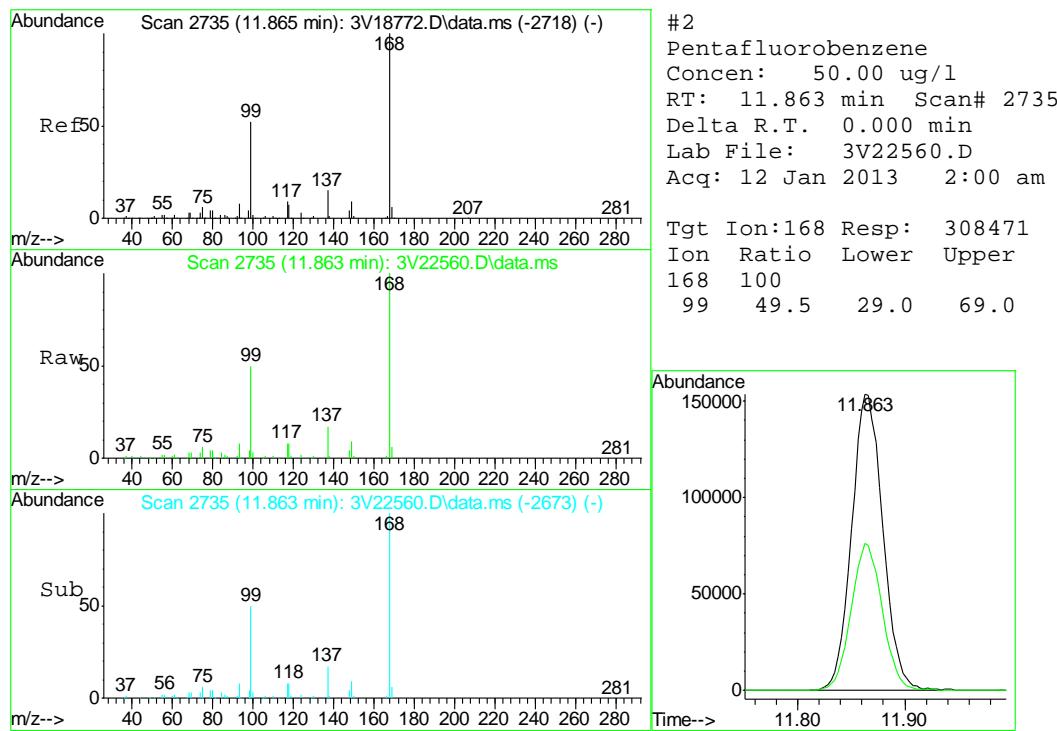
(#) = qualifier out of range (m) = manual integration (+) = signals summed

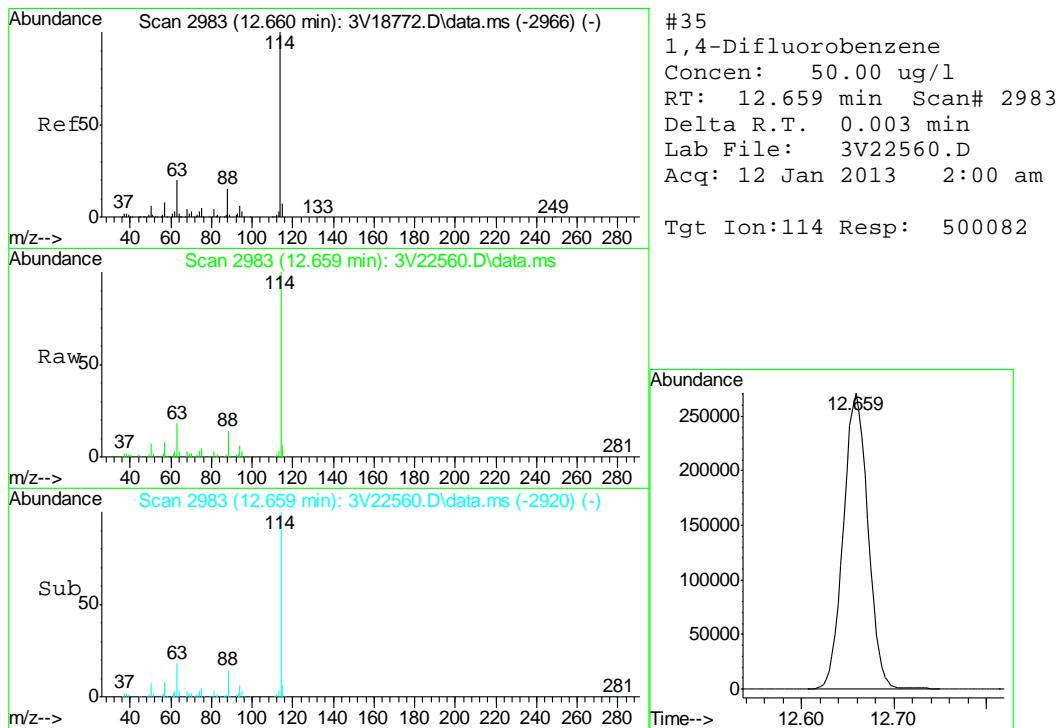
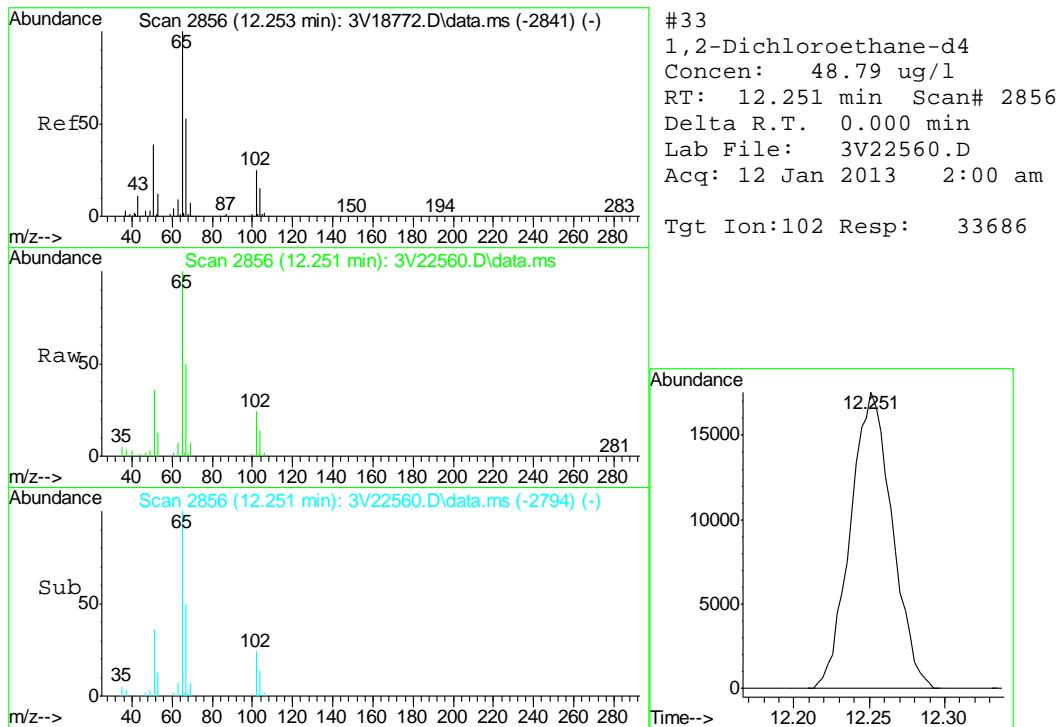
## Quantitation Report (QT Reviewed)

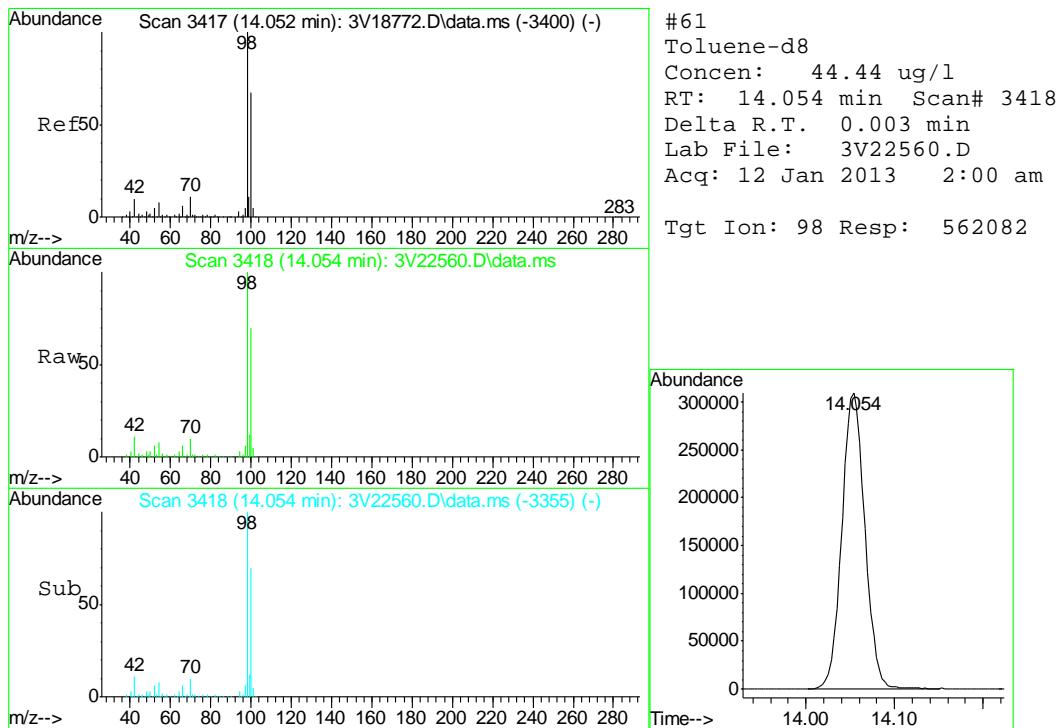
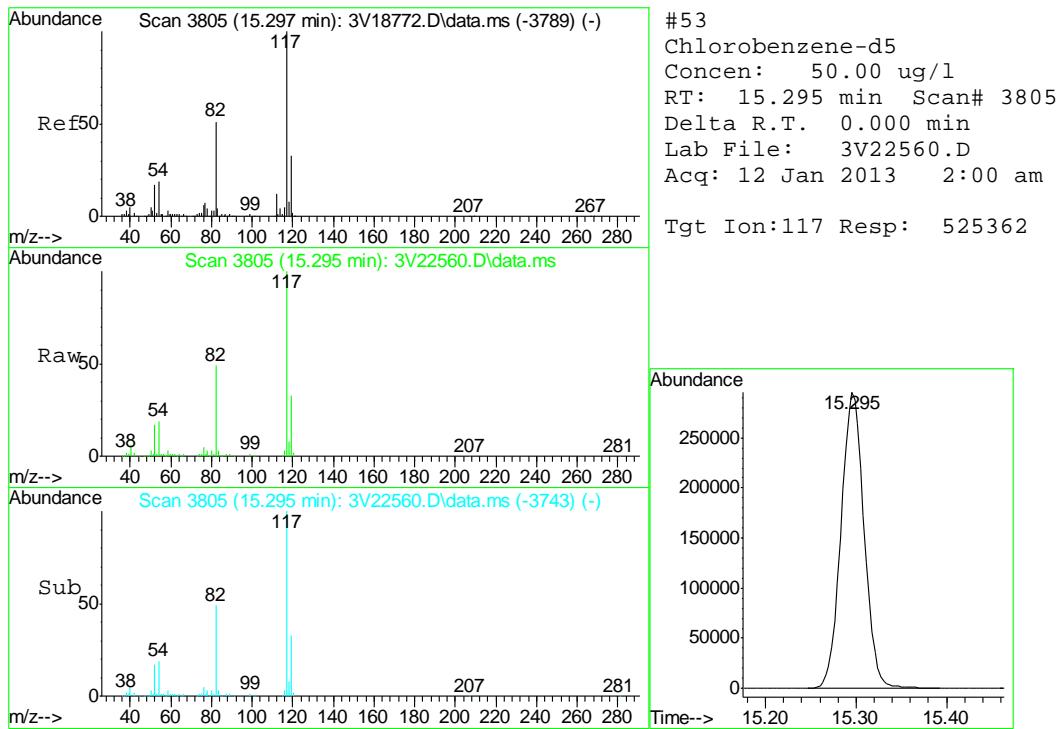
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 Data File : 3V22560.D  
 Acq On : 12 Jan 2013 2:00 am  
 Operator : BRETD  
 Sample : MB  
 Misc : MS5218,V3V1327,5.00,,100,5,1  
 ALS Vial : 27 Sample Multiplier: 1

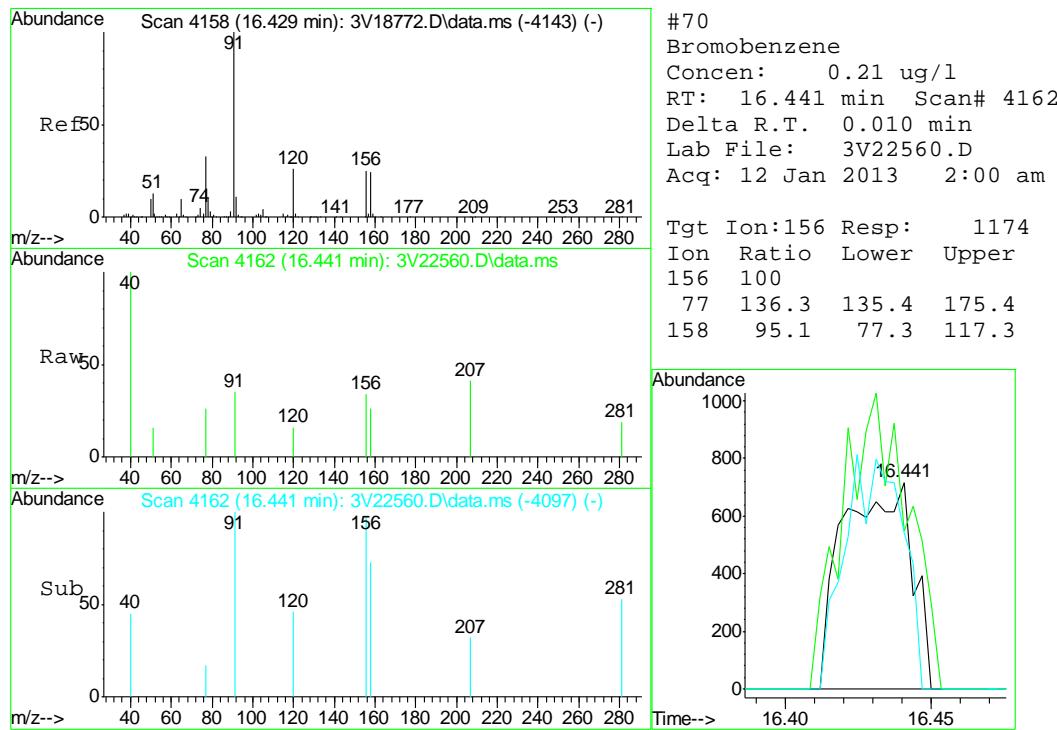
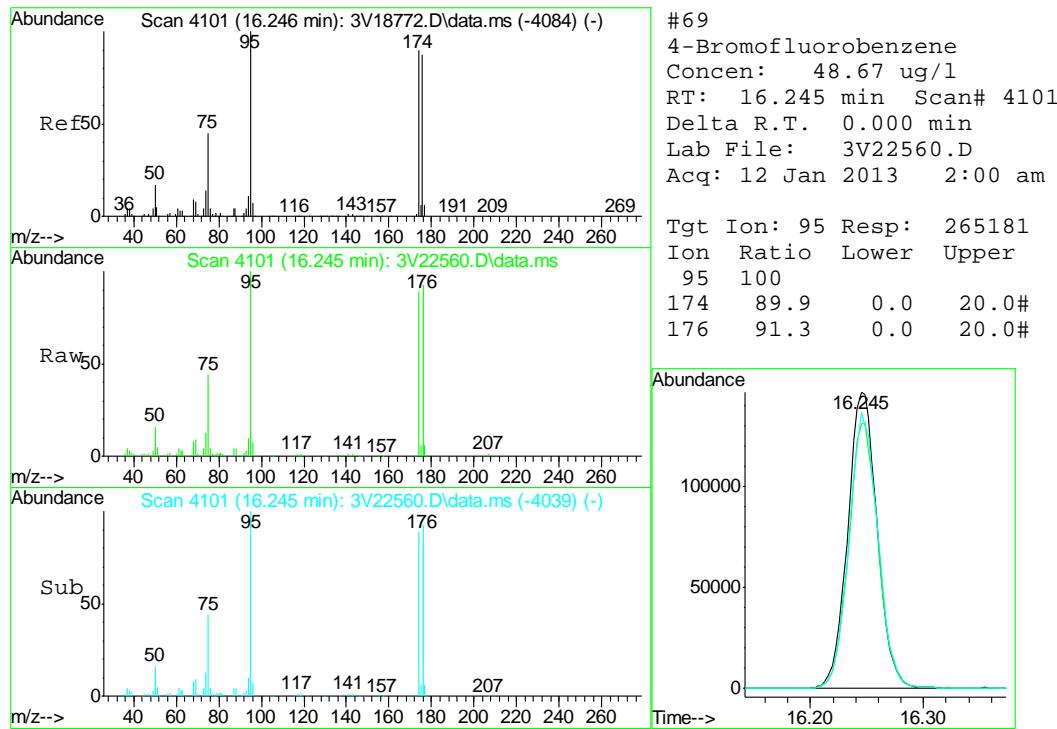
Quant Time: Jan 15 08:11:29 2013  
 Quant Method : C:\msdchem\1\METHODS\V3AP1299TVH1299SOIL.M  
 Quant Title : 8260  
 QLast Update : Thu Jan 03 11:40:16 2013  
 Response via : Initial Calibration

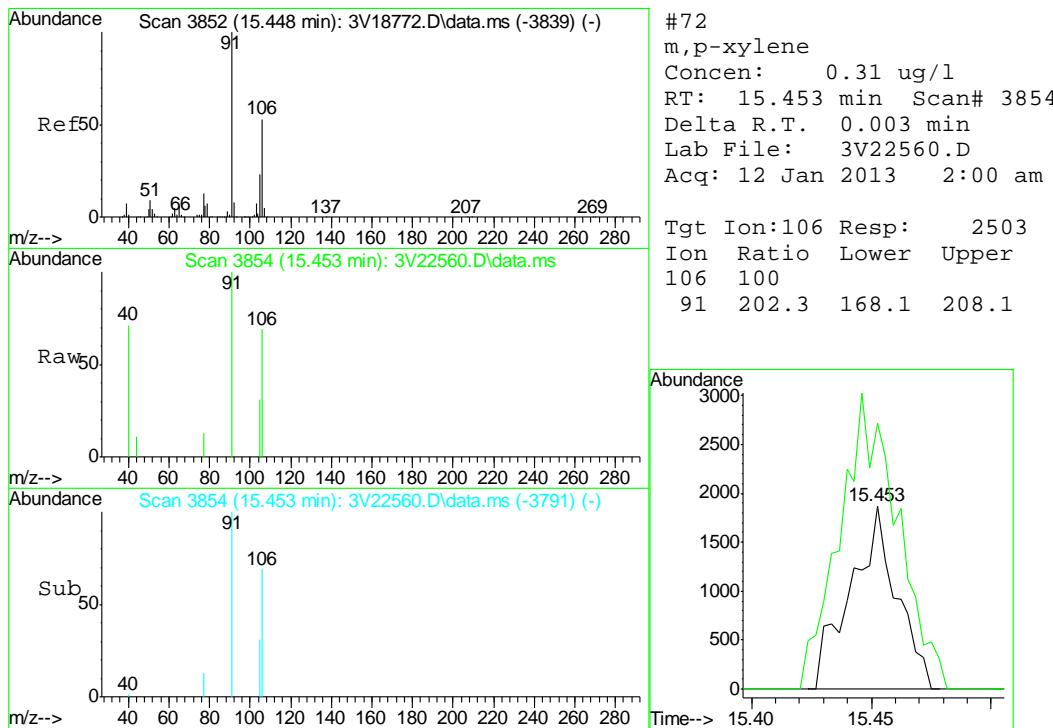
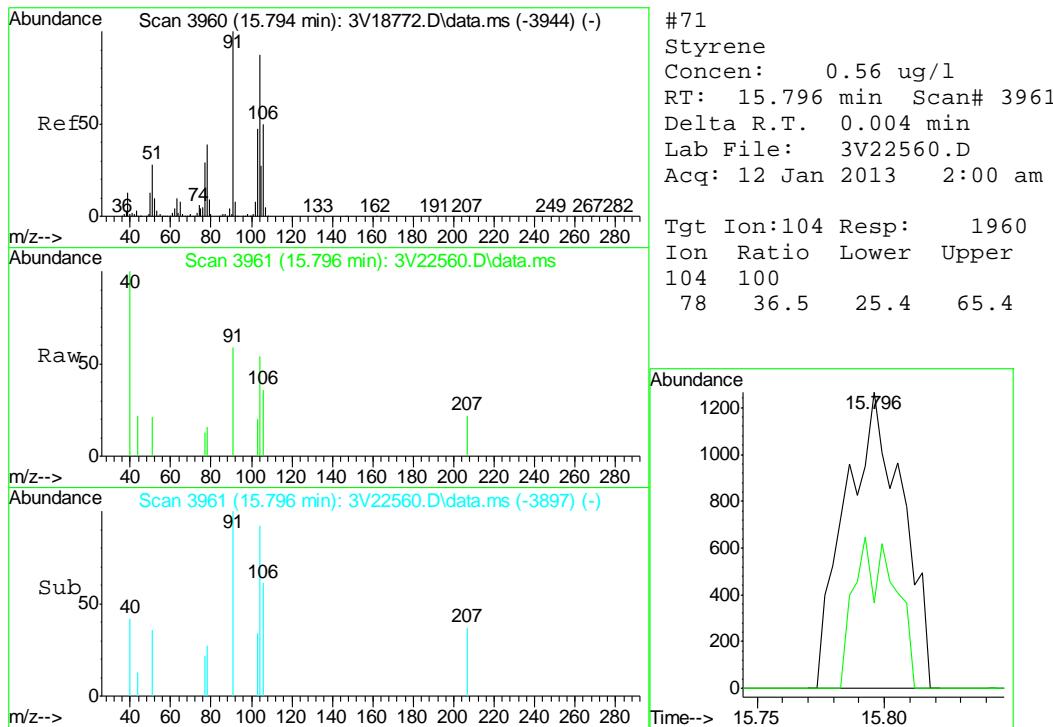


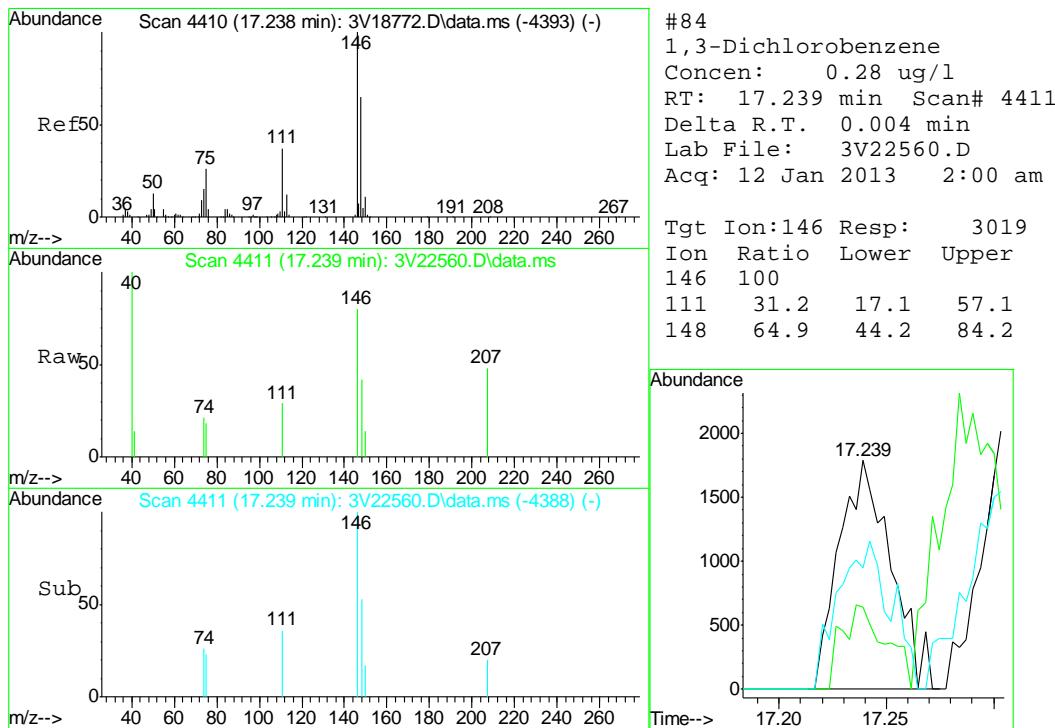
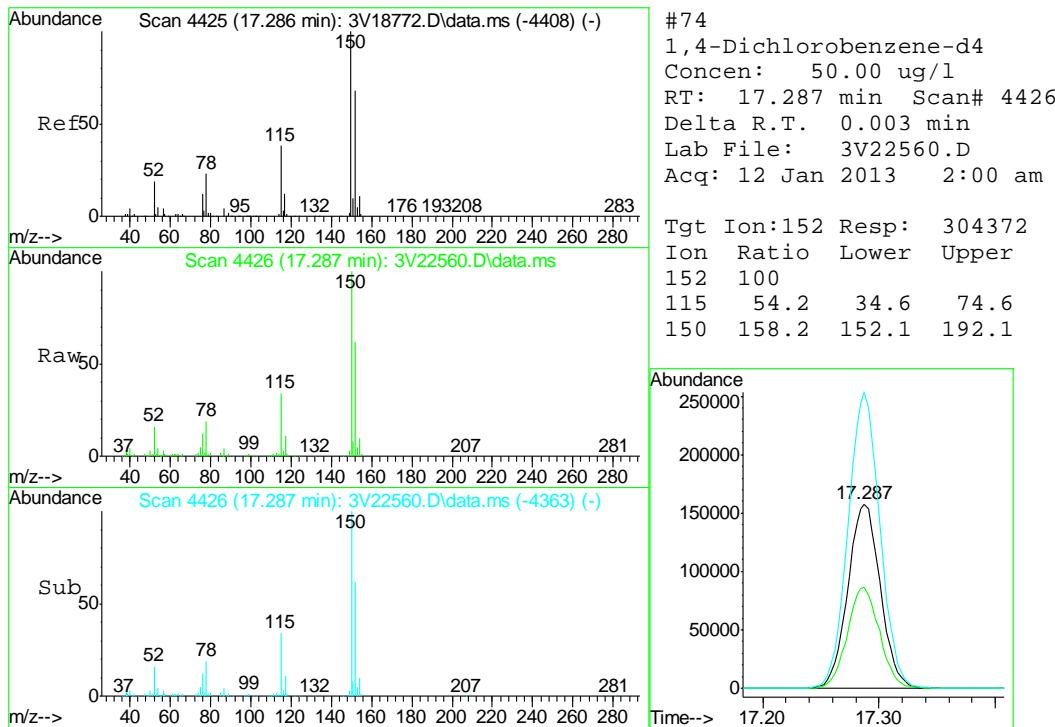


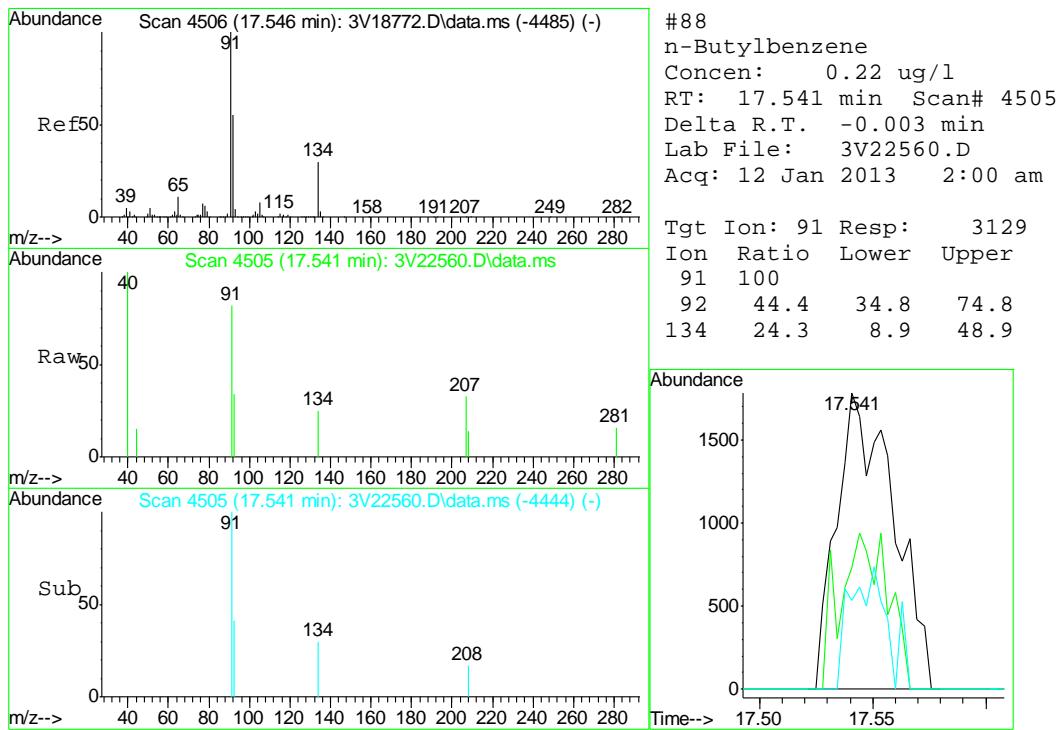
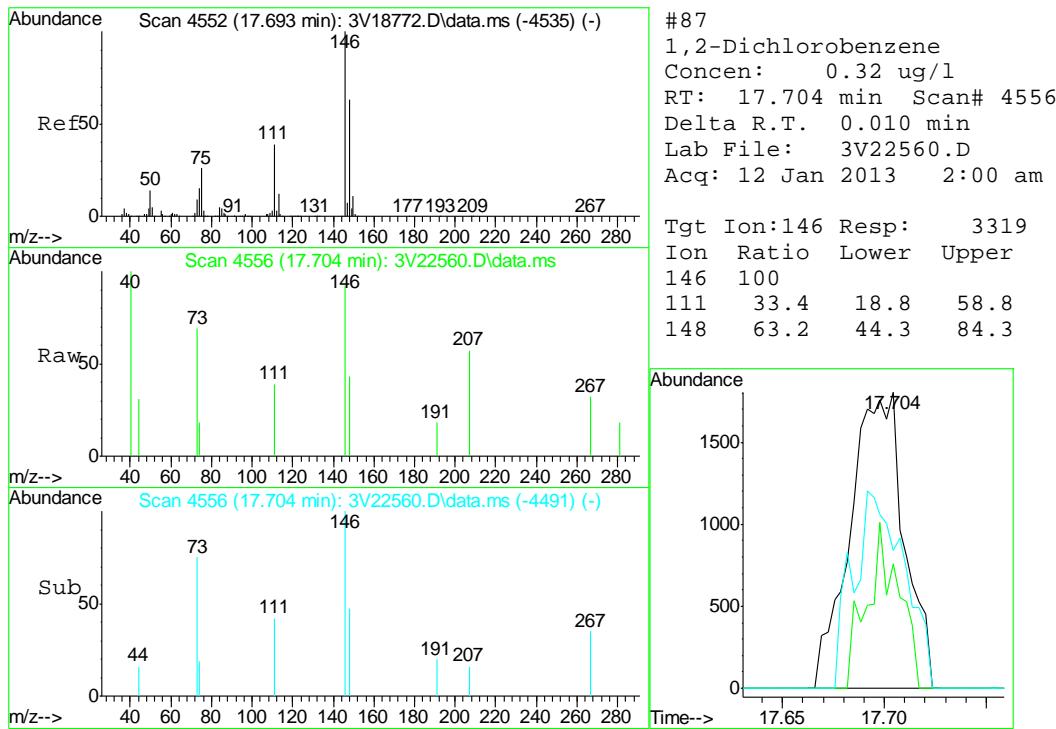


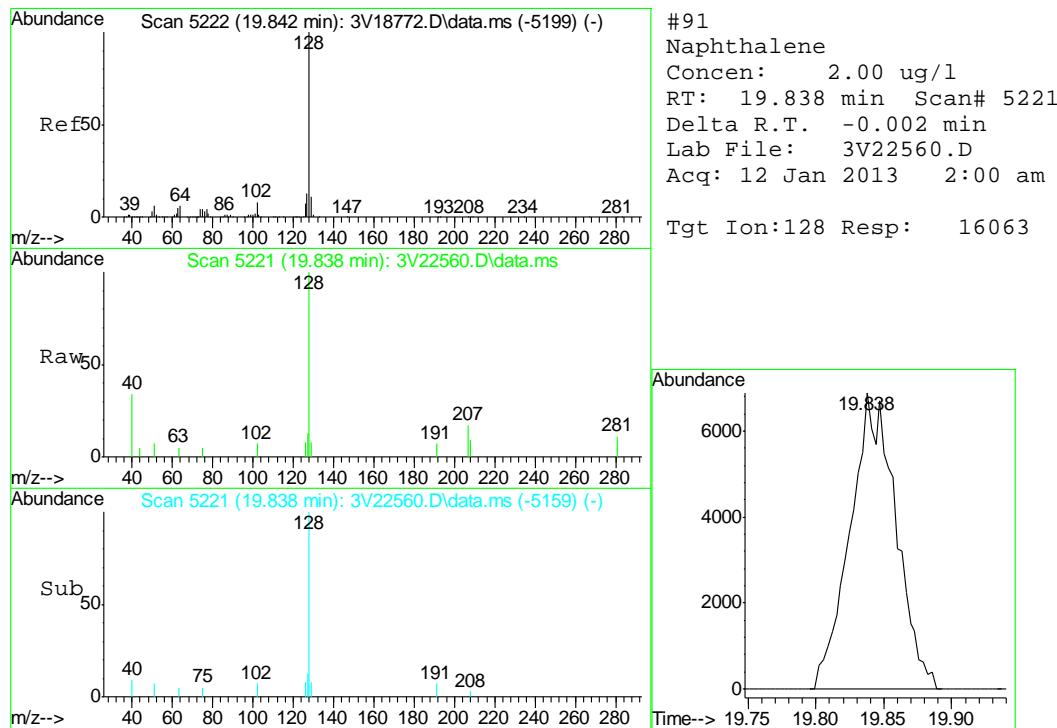
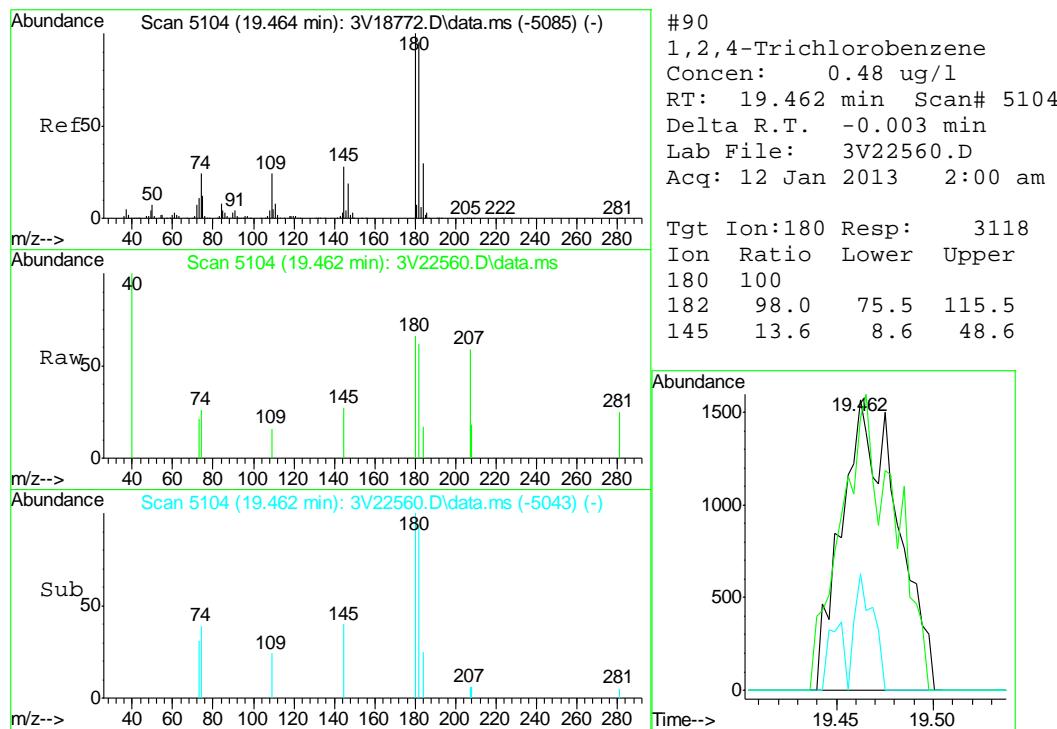


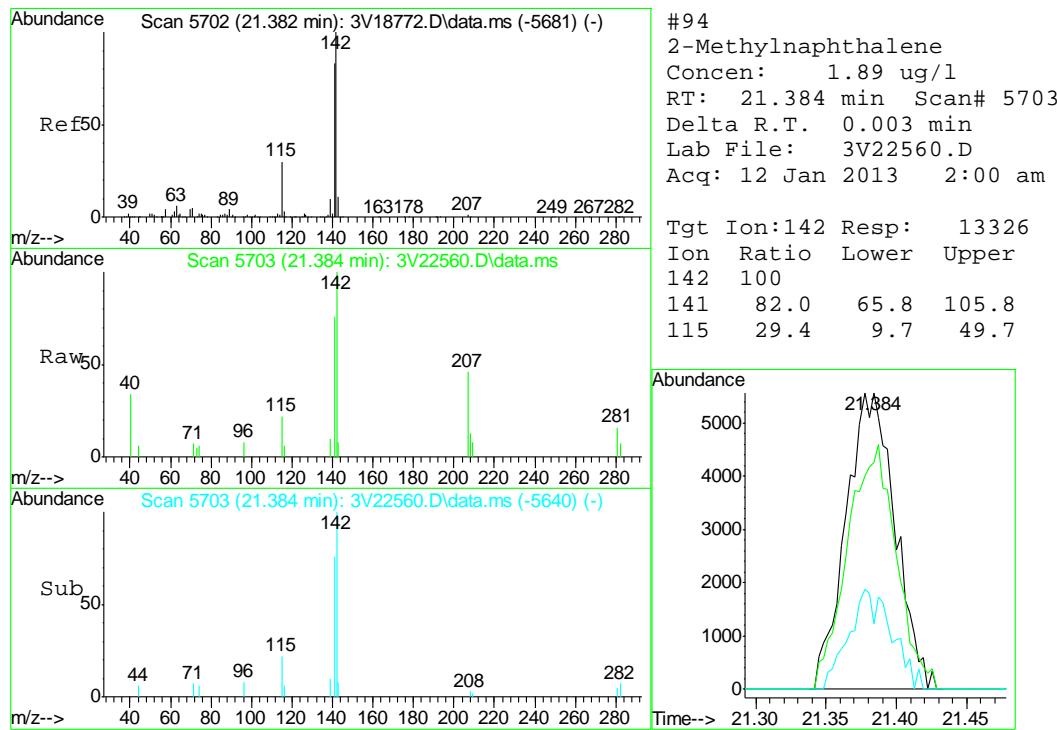
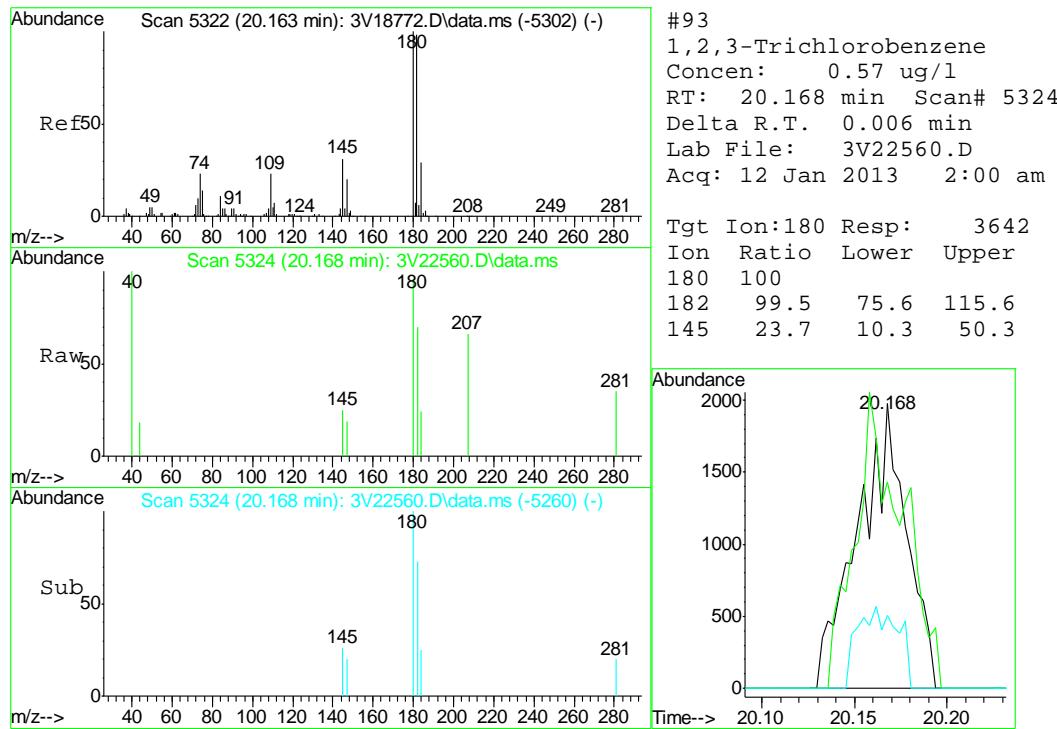


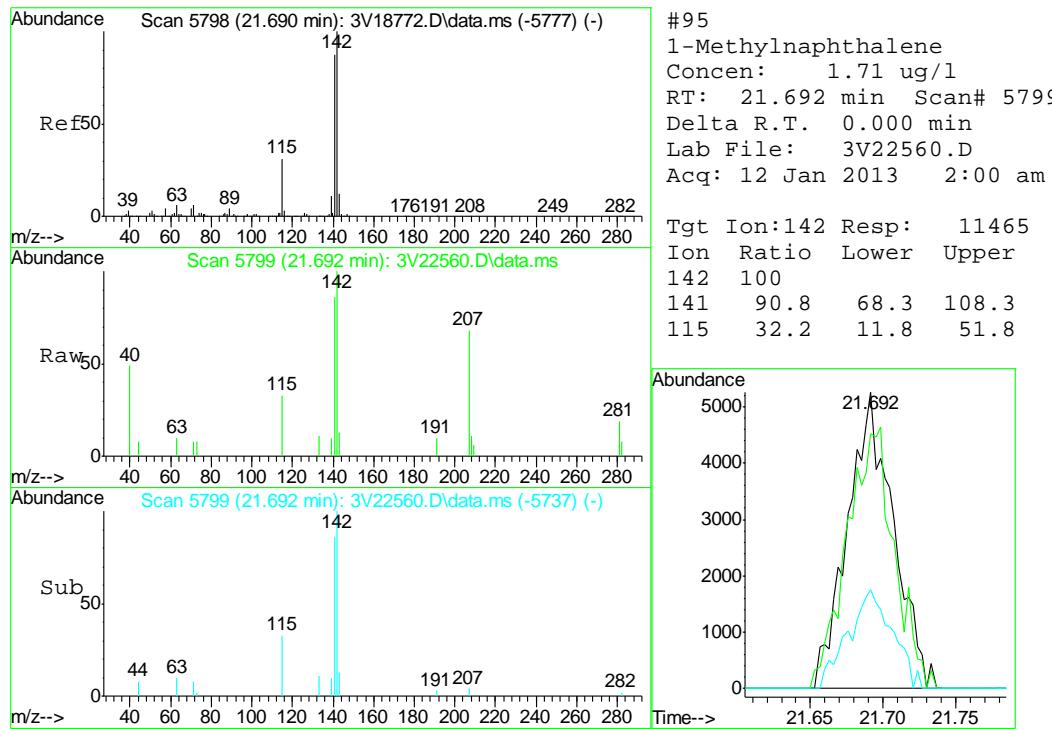














## GC/MS Semi-volatiles

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### QC Data Summaries

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Includes the following where applicable:

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

**Method Blank Summary**

**Job Number:** D42556  
**Account:** XTOKWR XTO Energy  
**Project:** PCU 296-5A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
OP7223-MB	3G12973.D	1	01/15/13	DC	01/14/13	OP7223	E3G621

The QC reported here applies to the following samples:

**Method:** SW846 8270C BY SIM

D42556-1, D42556-2

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	81% 10-159%
321-60-8	2-Fluorobiphenyl	85% 19-131%
1718-51-0	Terphenyl-d14	105% 18-150%

## Blank Spike Summary

Page 1 of 1

Job Number: D42556

Account: XTOKWR XTO Energy

Project: PCU 296-5A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
OP7223-BS	3G12974.D	1	01/15/13	DC	01/14/13	OP7223	E3G621

The QC reported here applies to the following samples:

Method: SW846 8270C BY SIM

D42556-1, D42556-2

CAS No.	Compound	Spike ug/kg	BSP ug/kg	BSP %	Limits
83-32-9	Acenaphthene	83.3	87.0	104	68-130
120-12-7	Anthracene	83.3	75.1	90	67-130
56-55-3	Benzo(a)anthracene	83.3	79.3	95	65-130
205-99-2	Benzo(b)fluoranthene	83.3	79.2	95	44-130
207-08-9	Benzo(k)fluoranthene	83.3	65.4	78	56-131
50-32-8	Benzo(a)pyrene	83.3	70.8	85	62-130
218-01-9	Chrysene	83.3	70.3	84	70-130
53-70-3	Dibenzo(a,h)anthracene	83.3	79.8	96	55-130
206-44-0	Fluoranthene	83.3	76.4	92	70-130
86-73-7	Fluorene	83.3	80.9	97	70-130
193-39-5	Indeno(1,2,3-cd)pyrene	83.3	77.2	93	56-130
91-20-3	Naphthalene	83.3	71.1	85	70-130
129-00-0	Pyrene	83.3	74.5	89	70-130

CAS No.	Surrogate Recoveries	BSP	Limits
4165-60-0	Nitrobenzene-d5	103%	10-159%
321-60-8	2-Fluorobiphenyl	97%	19-131%
1718-51-0	Terphenyl-d14	107%	18-150%

\* = Outside of Control Limits.

# Matrix Spike/Matrix Spike Duplicate Summary

Page 1 of 1

Job Number: D42556

Account: XTOKWR XTO Energy

Project: PCU 296-5A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
OP7223-MS	3G12984.D	1	01/15/13	DC	01/14/13	OP7223	E3G621
OP7223-MSD	3G12985.D	1	01/15/13	DC	01/14/13	OP7223	E3G621
D42510-1	3G12983.D	1	01/15/13	DC	01/14/13	OP7223	E3G621

The QC reported here applies to the following samples:

Method: SW846 8270C BY SIM

D42556-1, D42556-2

CAS No.	Compound	D42510-1		Spike	MS	MS	MSD	MSD	RPD	Limits Rec/RPD
		ug/kg	Q	ug/kg	ug/kg	%	ug/kg	%		
83-32-9	Acenaphthene	ND		98.3	76.7	78	76.8	78	0	25-151/30
120-12-7	Anthracene	ND		98.3	93.2	95	95.1	97	2	39-159/30
56-55-3	Benzo(a)anthracene	ND		98.3	107	109	109	111	2	39-168/30
205-99-2	Benzo(b)fluoranthene	ND		98.3	109	111	110	112	1	24-163/30
207-08-9	Benzo(k)fluoranthene	ND		98.3	70.7	72	74.9	76	6	10-188/30
50-32-8	Benzo(a)pyrene	ND		98.3	77.1	78	79.0	80	2	32-144/30
218-01-9	Chrysene	25.4		98.3	97.1	73	101	77	4	43-150/30
53-70-3	Dibenzo(a,h)anthracene	ND		98.3	76.5	78	79.8	81	4	21-152/30
206-44-0	Fluoranthene	ND		98.3	101	103	102	104	1	36-157/30
86-73-7	Fluorene	198		98.3	260	63	249	52	4	10-182/30
193-39-5	Indeno(1,2,3-cd)pyrene	ND		98.3	73.2	74	76.2	77	4	20-154/30
91-20-3	Naphthalene	1070		98.3	1400	336* a	909	-164* a	43* b	10-163/30
129-00-0	Pyrene	31.3		98.3	126	96	127	97	1	25-180/30

CAS No.	Surrogate Recoveries	MS	MSD	D42510-1	Limits
4165-60-0	Nitrobenzene-d5	141%	110%	141%	10-159%
321-60-8	2-Fluorobiphenyl	63%	56%	61%	19-131%
1718-51-0	Terphenyl-d14	98%	96%	91%	18-150%

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

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

\* = Outside of Control Limits.

8.3.1  
8



## GC/MS Semi-volatiles

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Raw Data

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Judy Nelson  
 01/16/13 11:51

## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\011513\  
 Data File : 3g12980.D  
 Acq On : 15 Jan 2013 1:06 pm  
 Operator : DONC  
 Sample : D42556-1  
 Misc : OP7223,E3G621,30.06,,,1,1  
 ALS Vial : 11 Sample Multiplier: 1

Quant Time: Jan 15 13:59:15 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	5.620	136	132711	4.0000	ug/mL	0.00
6) Acenaphthene-d10	7.337	164	78214	4.0000	ug/mL	0.01
15) Phenanthrene-d10	8.812	188	131819	4.0000	ug/mL	0.00
19) Chrysene-d12	11.443	240	97134	4.0000	ug/mL	0.00
24) Perylene-d12	12.810	264	80773	4.0000	ug/mL	0.00

## System Monitoring Compounds

2) Nitrobenzene-d5	4.935	82	495666	41.5235	ug/mL	-0.01
Spiked Amount	50.000	Range	25 - 135	Recovery	=	83.04%
7) 2-Fluorobiphenyl	6.664	172	1296043	43.4882	ug/mL	0.00
Spiked Amount	50.000	Range	25 - 135	Recovery	=	86.98%
21) Terphenyl-d14	10.402	244	714045	54.0249	ug/mL	0.00
Spiked Amount	50.000	Range	25 - 135	Recovery	=	108.04%

## Target Compounds

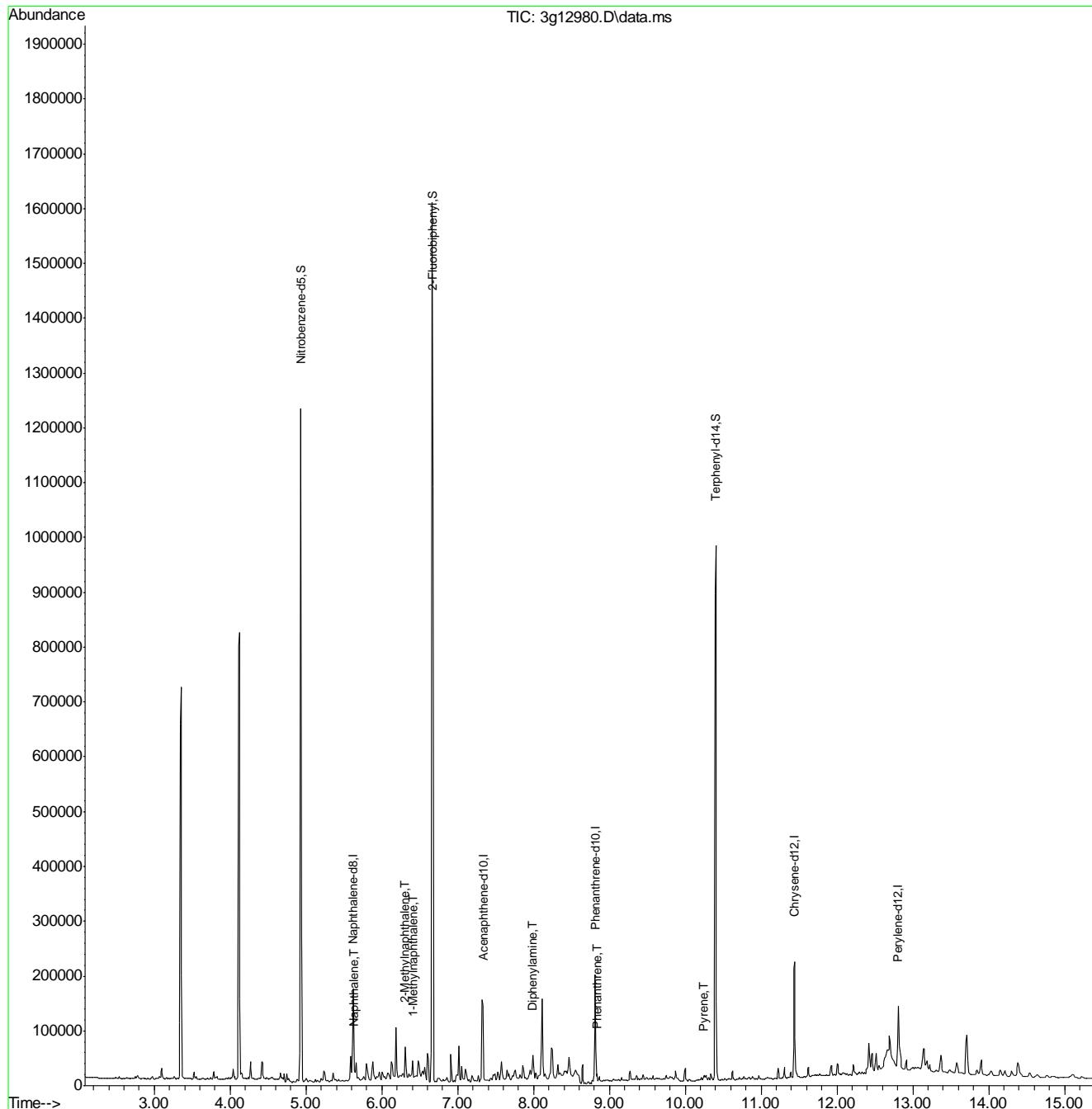
				Qvalue
3) N-Nitrosodimethylamine	2.312	74	50	N.D.
4) N-Nitrosodi-propylamine	0.000	70	0	N.D. d
5) Naphthalene	5.633	128	10781	0.2694 ug/mL# 50
8) 2-Methylnaphthalene	6.306	142	23572	0.9379 ug/mL 90
9) 1-Methylnaphthalene	6.406	142	11401	0.5186 ug/mL 88
10) Acenaphthylene	7.184	152	825	N.D.
11) Acenaphthene	7.361	154	327	N.D.
12) Dibenzofuran	0.000	168	0	N.D. d
13) Fluorene	7.869	166	1274	N.D.
14) Diphenylamine	7.987	169	24537m	0.9998 ug/mL
16) Phenanthrene	8.835	178	12035	0.2363 ug/mL# 60
17) Anthracene	8.922	178	452	N.D.
18) Fluoranthene	10.015	202	986	N.D.
20) Pyrene	10.244	202	5583	0.1076 ug/mL# 61
22) Benzo(a)anthracene	11.437	228	1261	N.D.
23) Chrysene	11.463	228	1173	N.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	14.408	276	469	N.D.

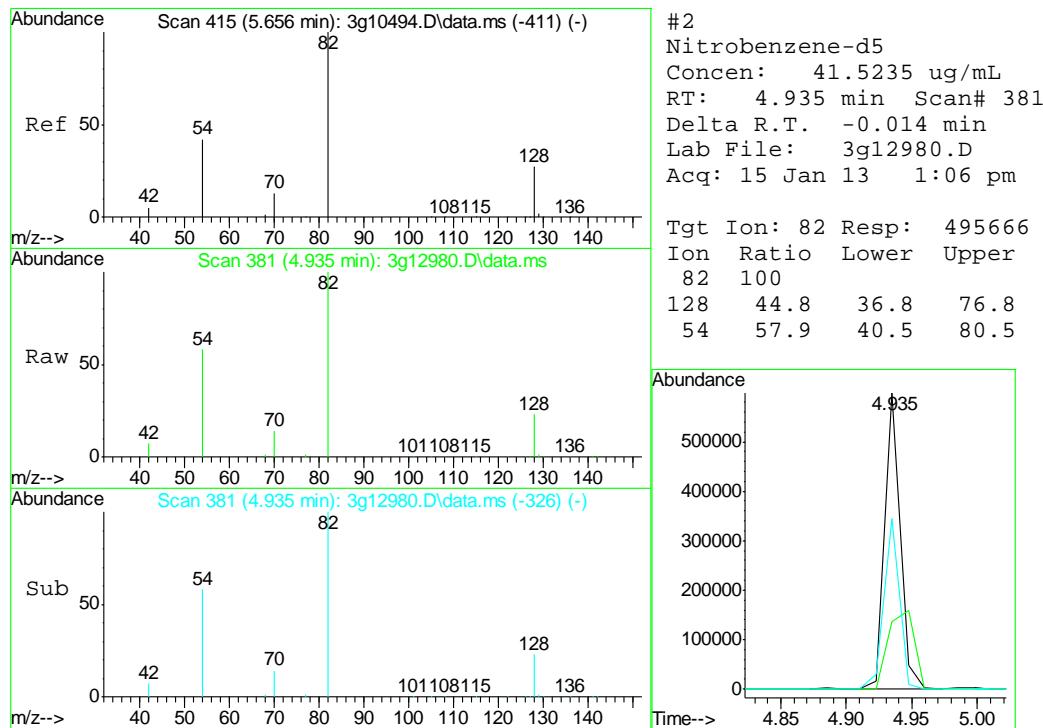
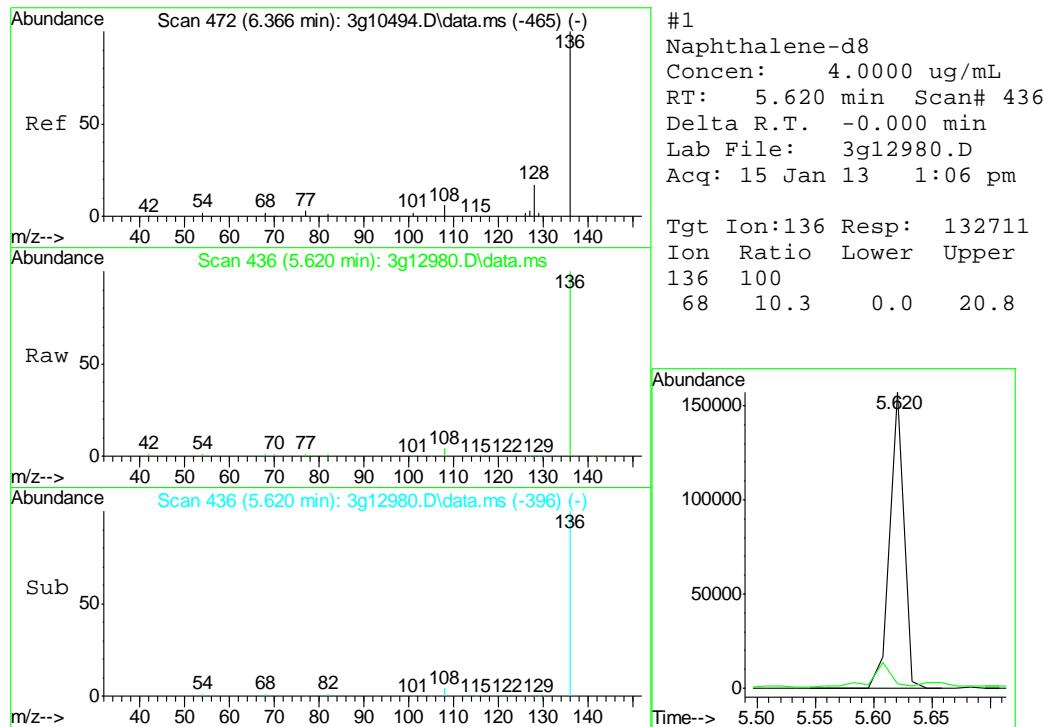
(#) = qualifier out of range (m) = manual integration (+) = signals summed

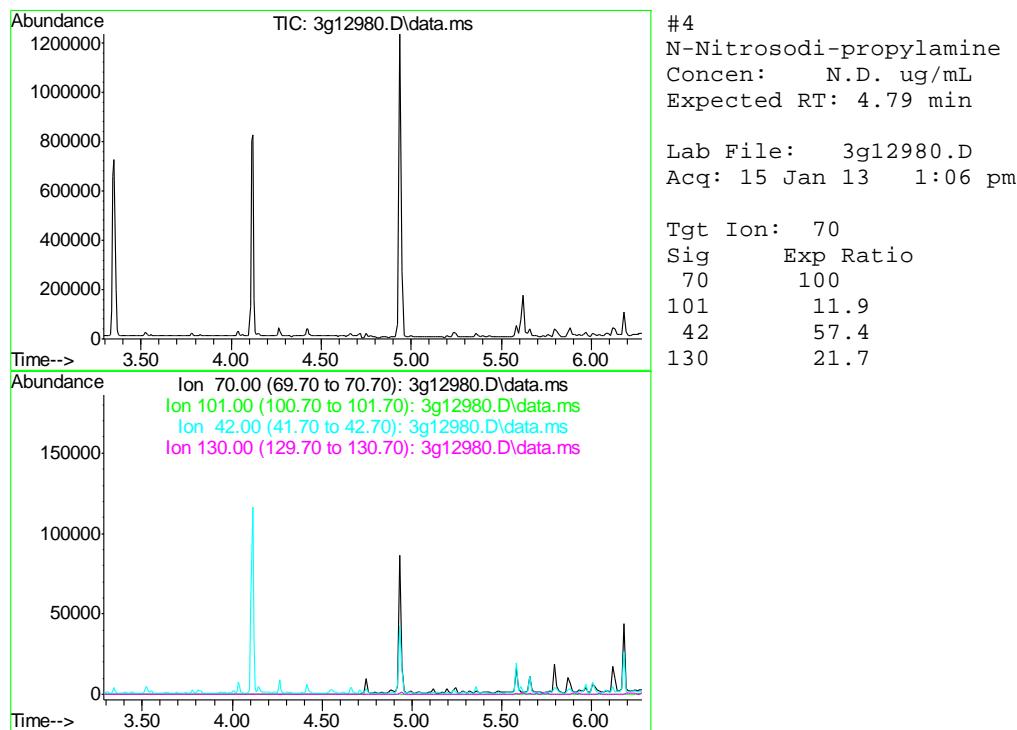
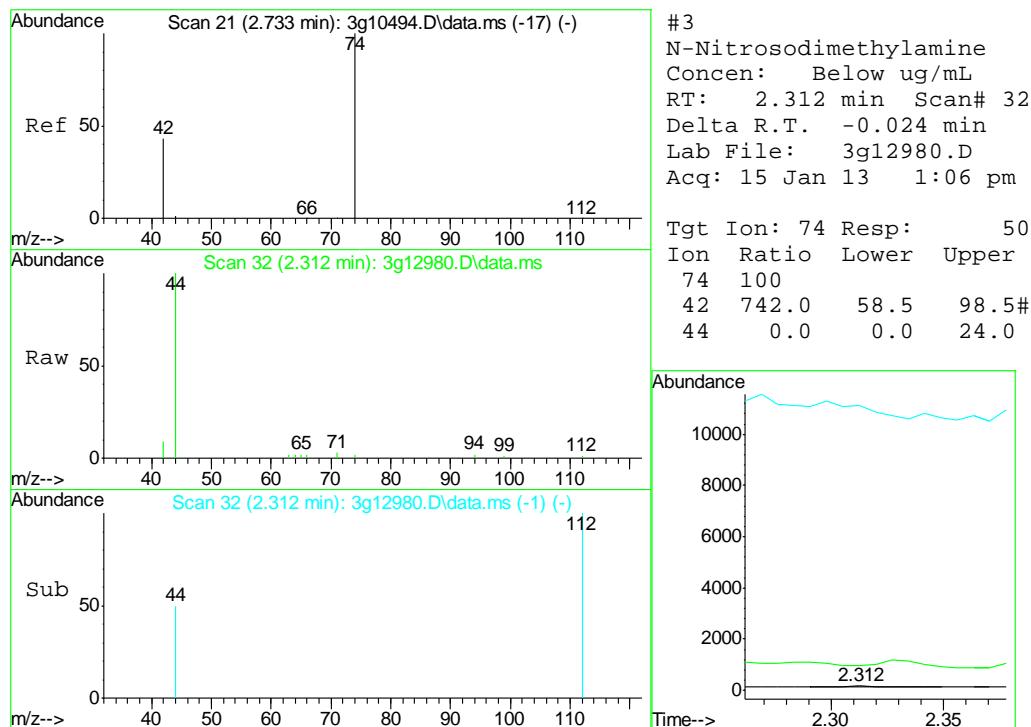
## Quantitation Report (QT Reviewed)

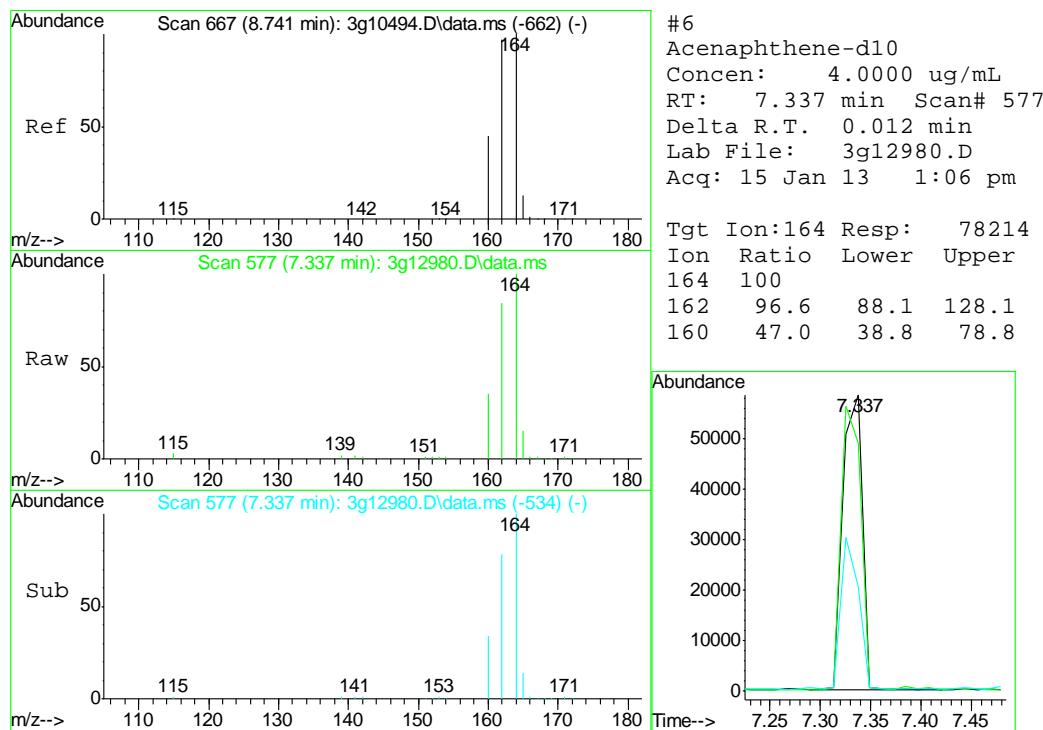
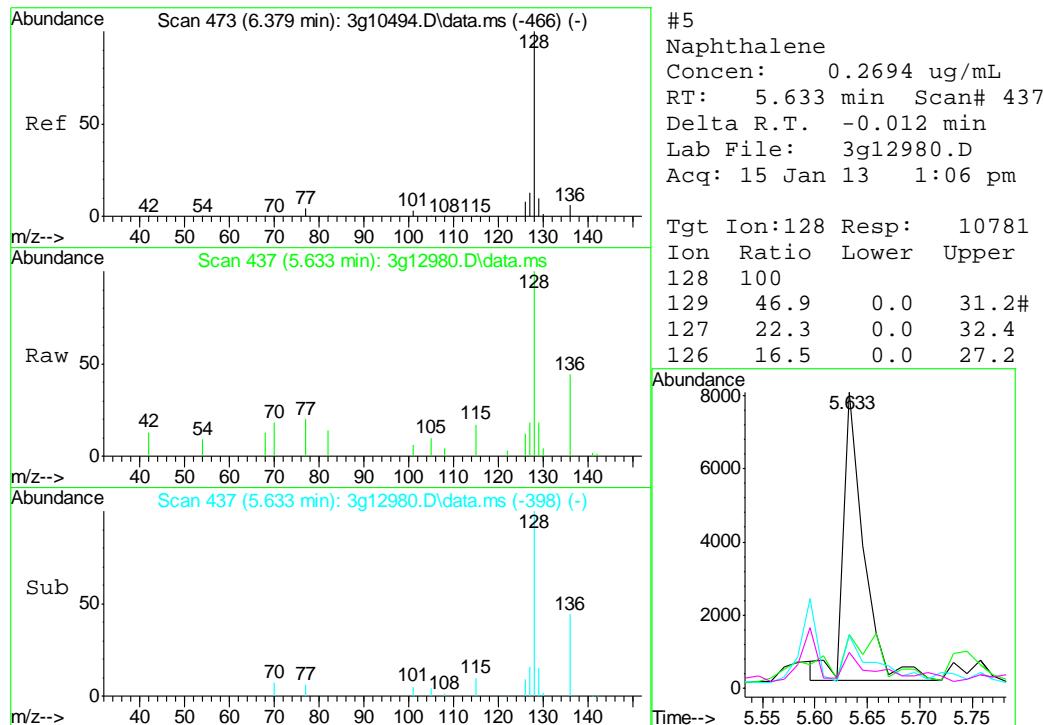
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 Data File : 3g12980.D  
 Acq On : 15 Jan 2013 1:06 pm  
 Operator : DONC  
 Sample : D42556-1  
 Misc : OP7223,E3G621,30.06,,,1,1  
 ALS Vial : 11 Sample Multiplier: 1

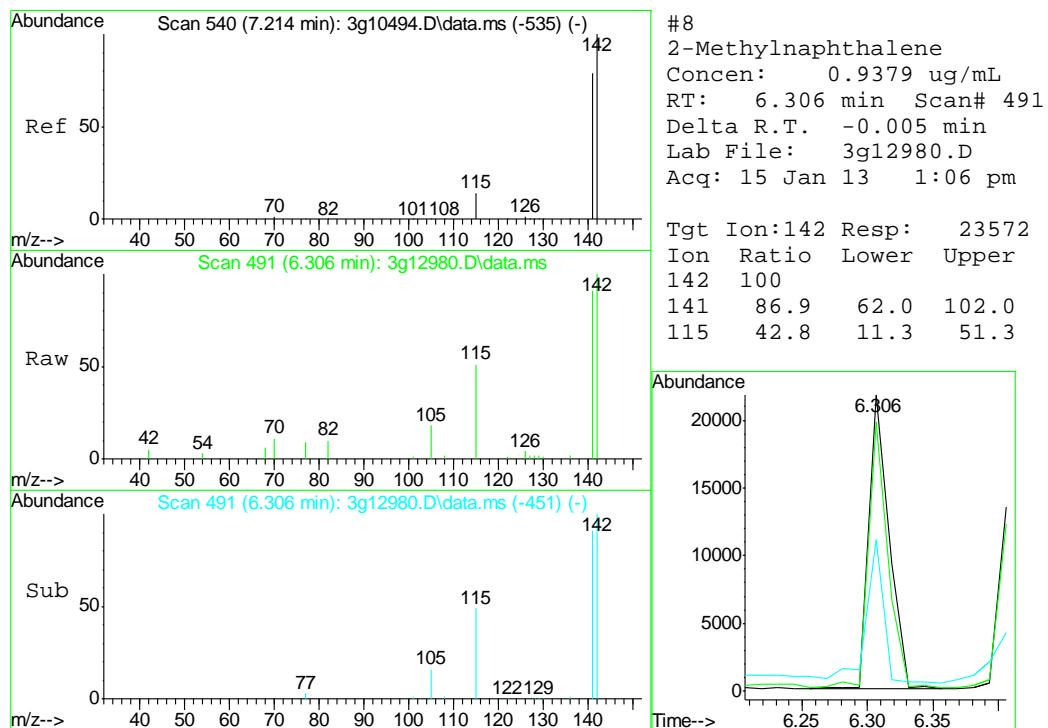
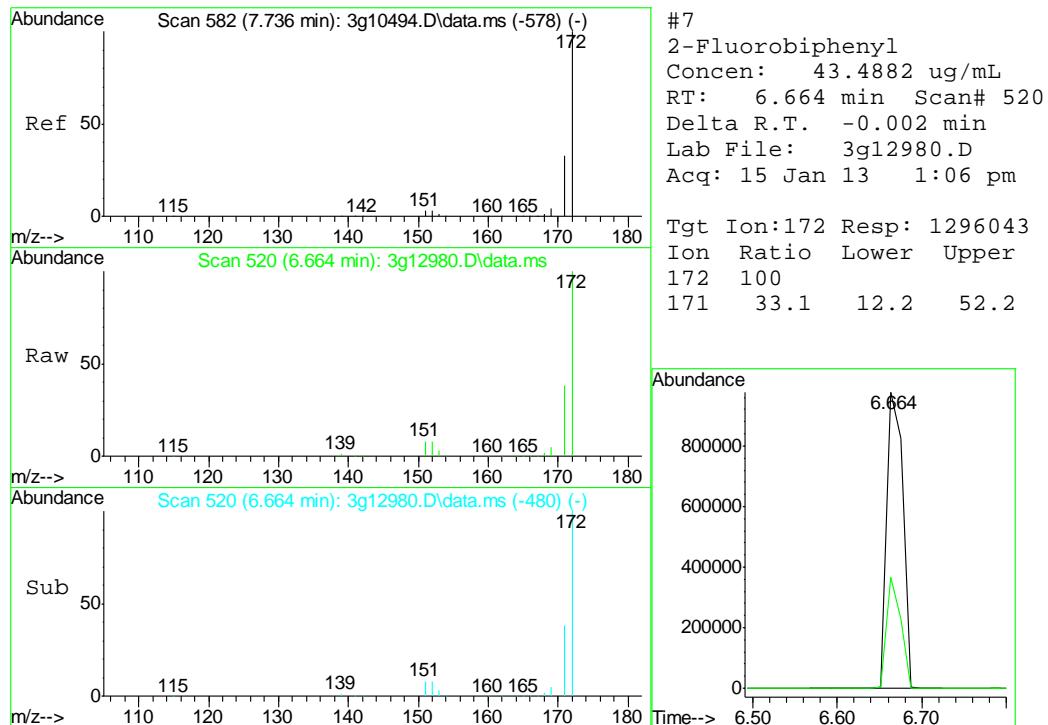
Quant Time: Jan 15 13:59:15 2013  
 Quant Method : C:\msdchem\1\METHODS\SIMPE3G611.M  
 Quant Title : PAHSIM BASE  
 QLast Update : Thu Jan 10 14:18:35 2013  
 Response via : Initial Calibration

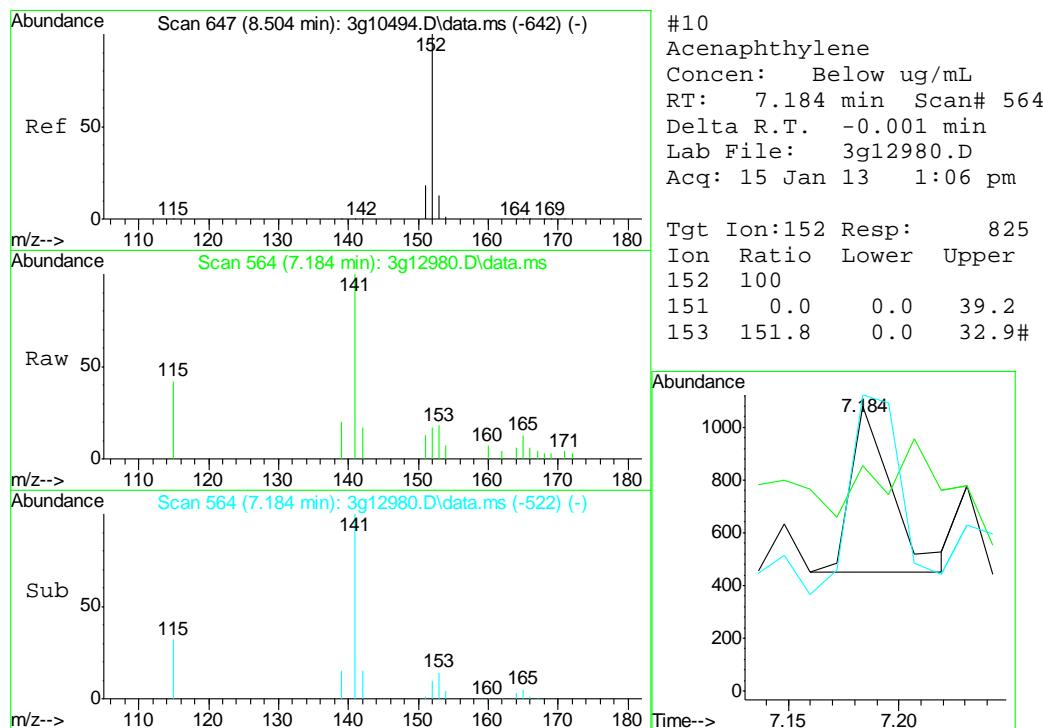
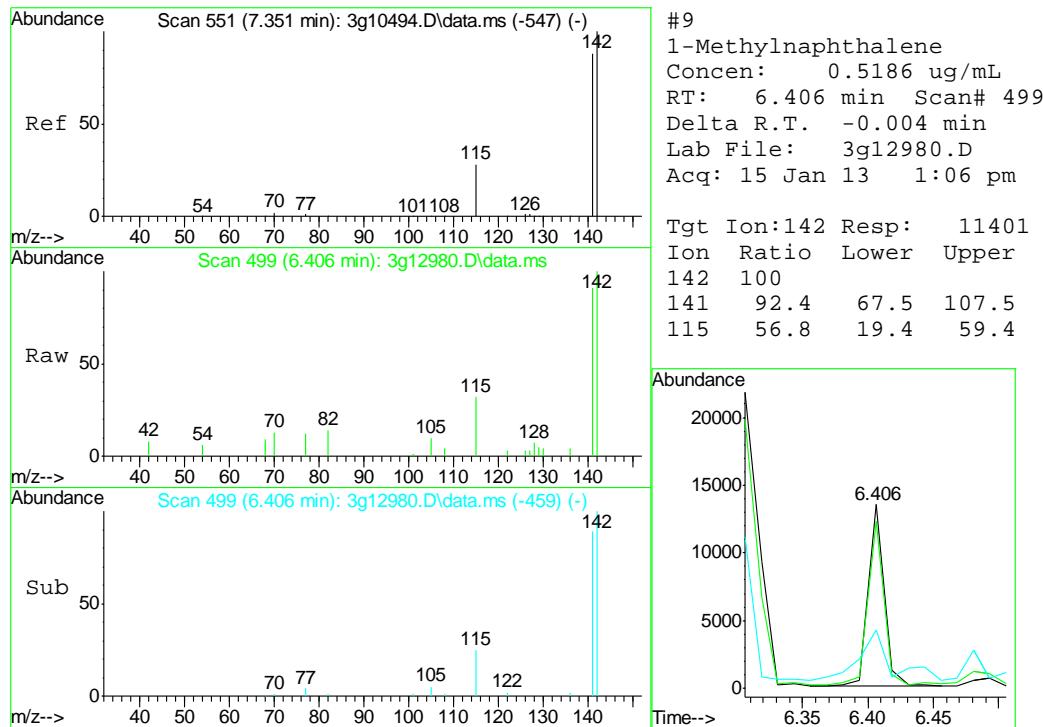


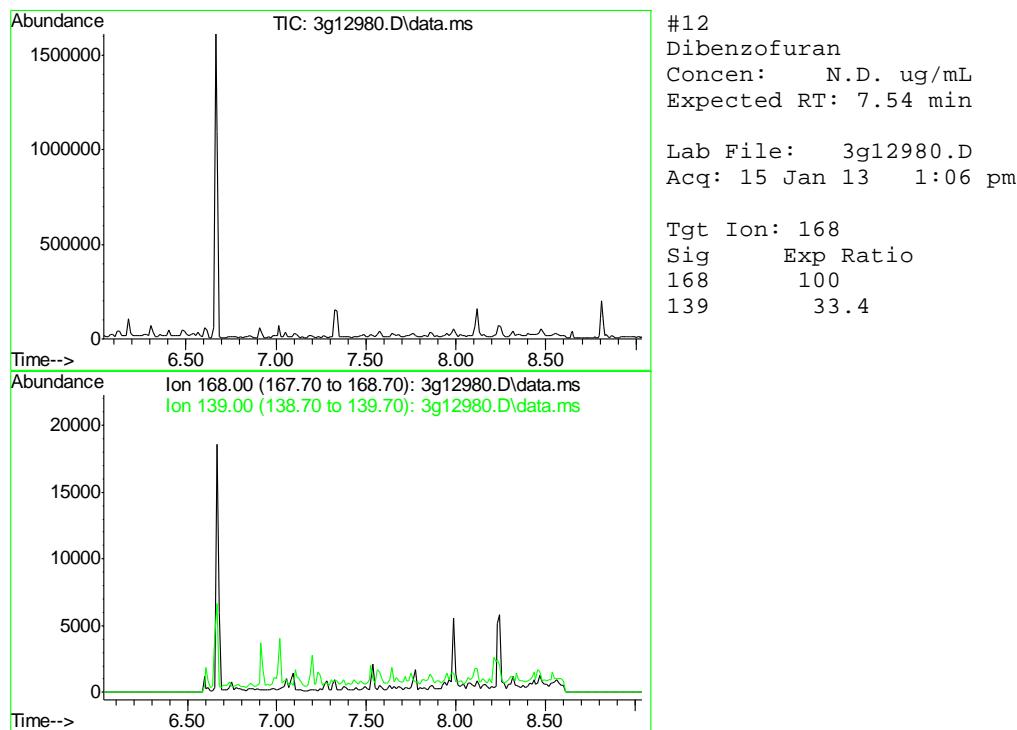
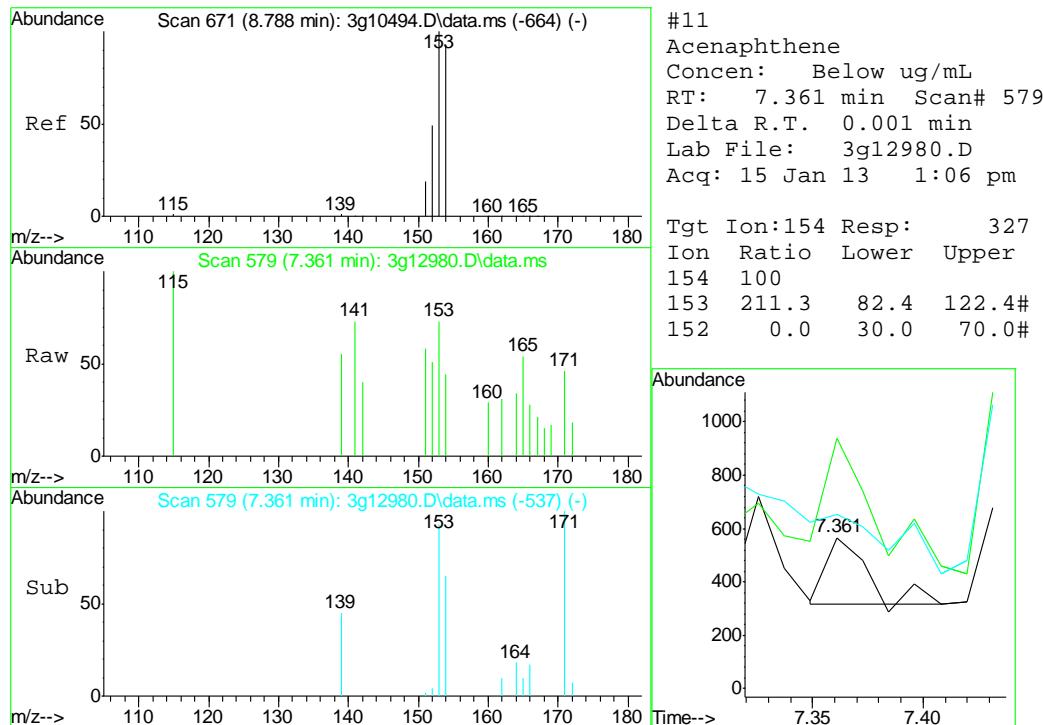


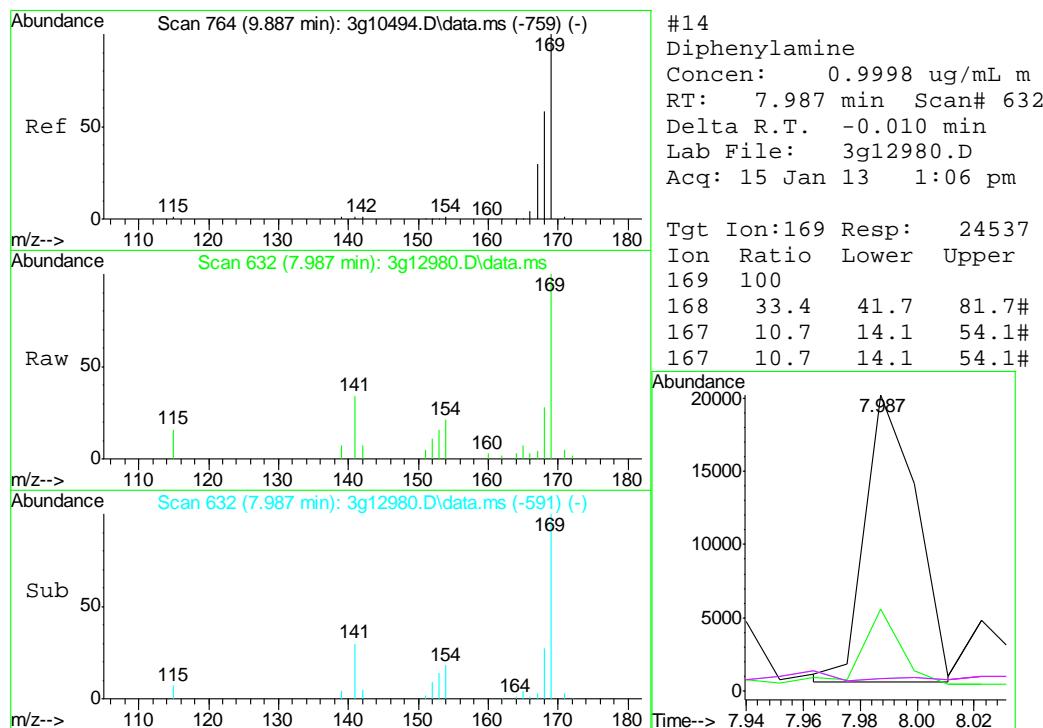
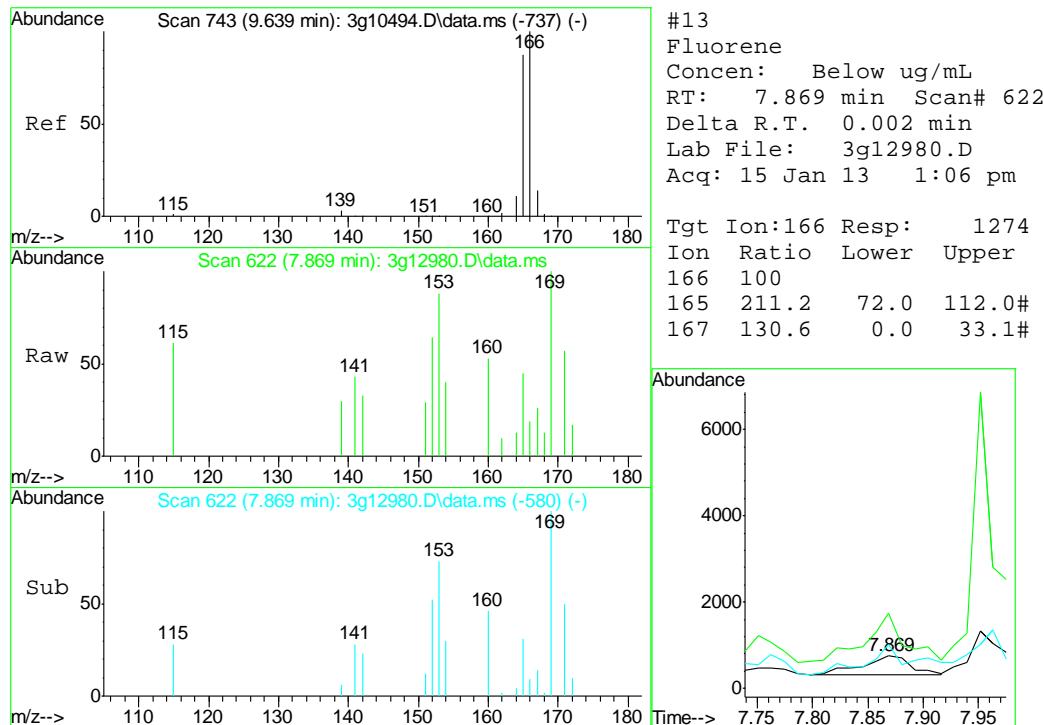


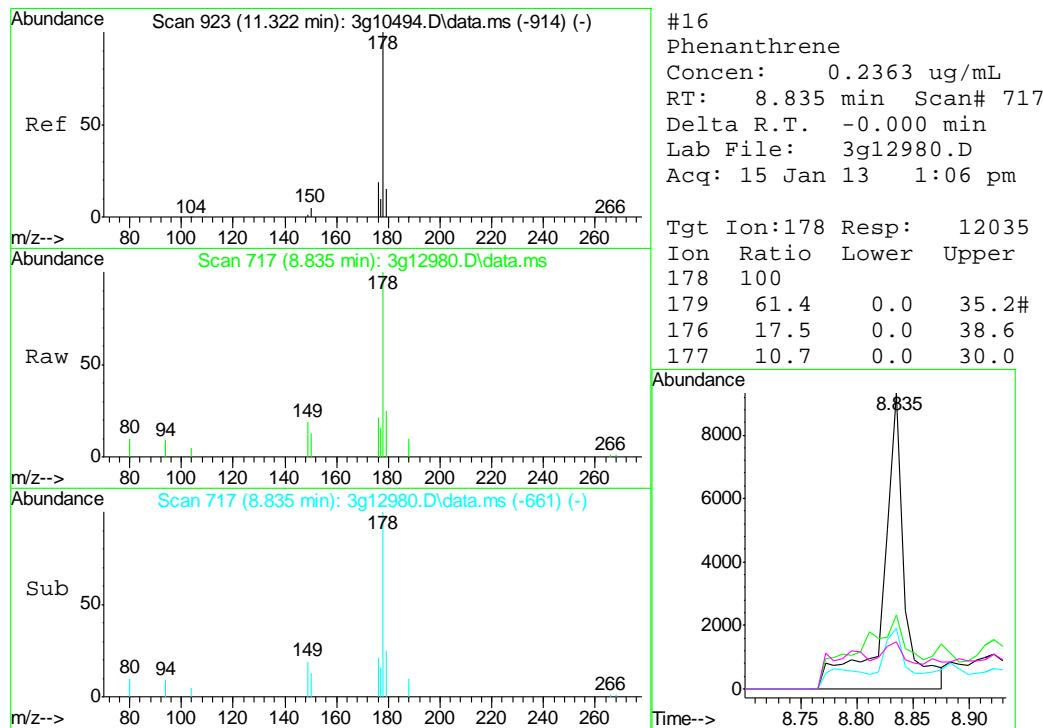
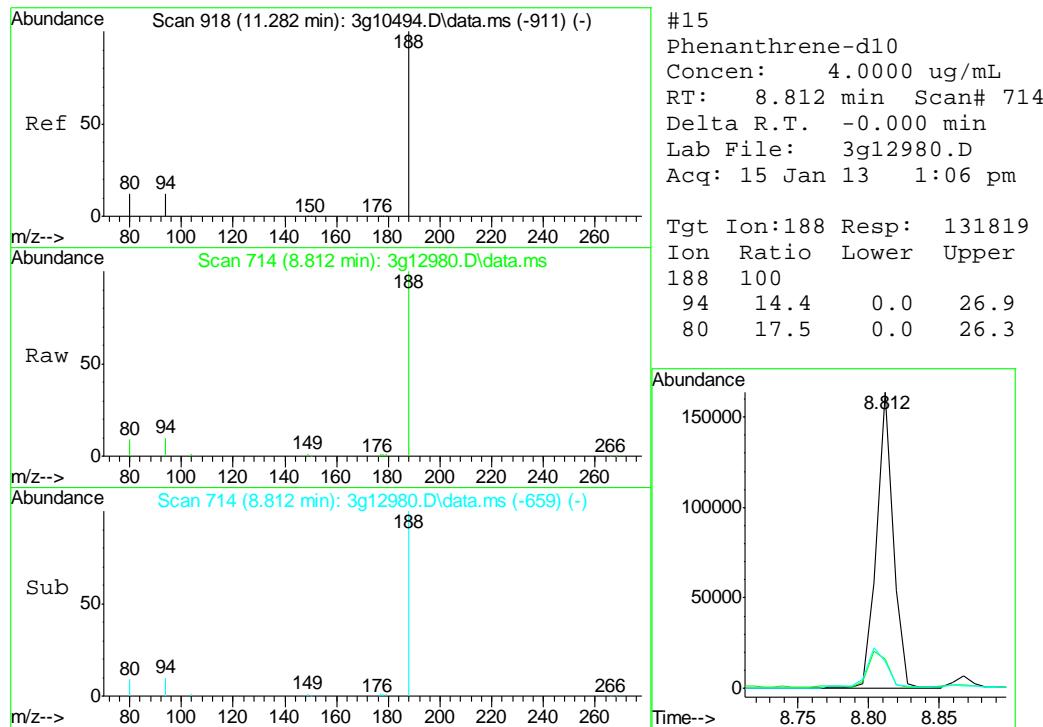


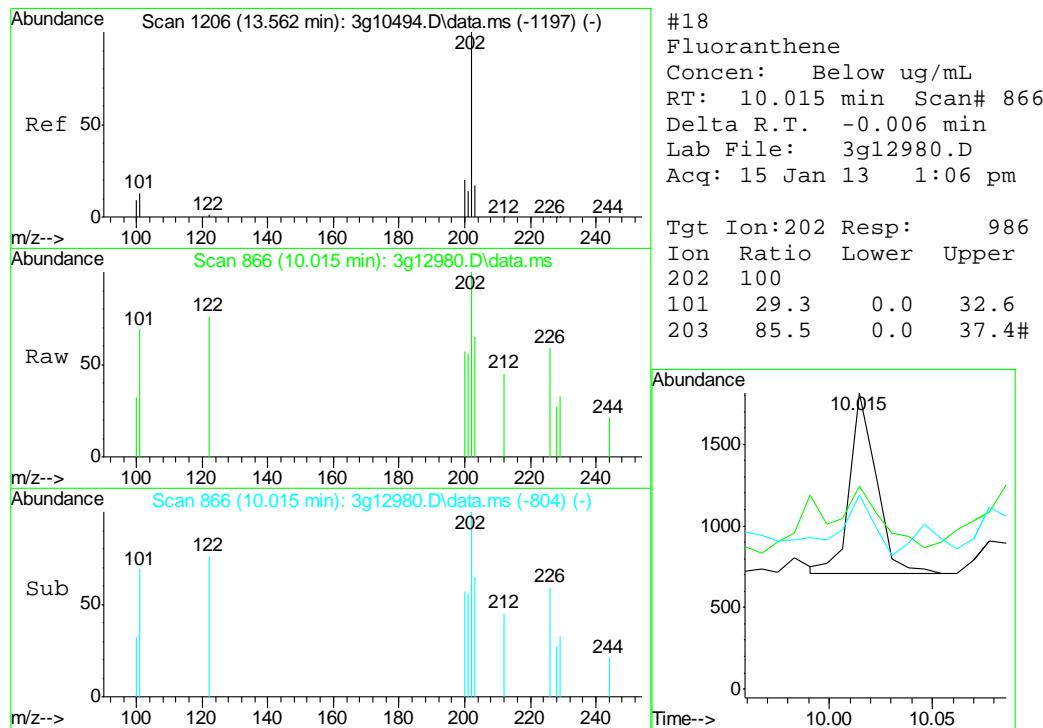
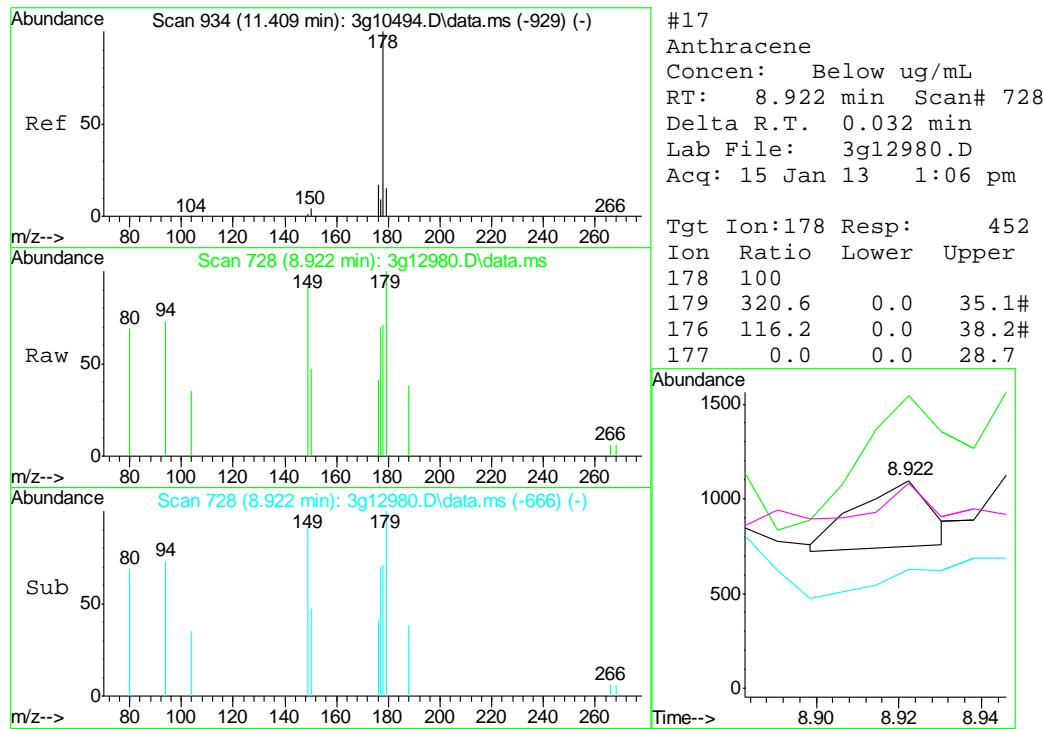


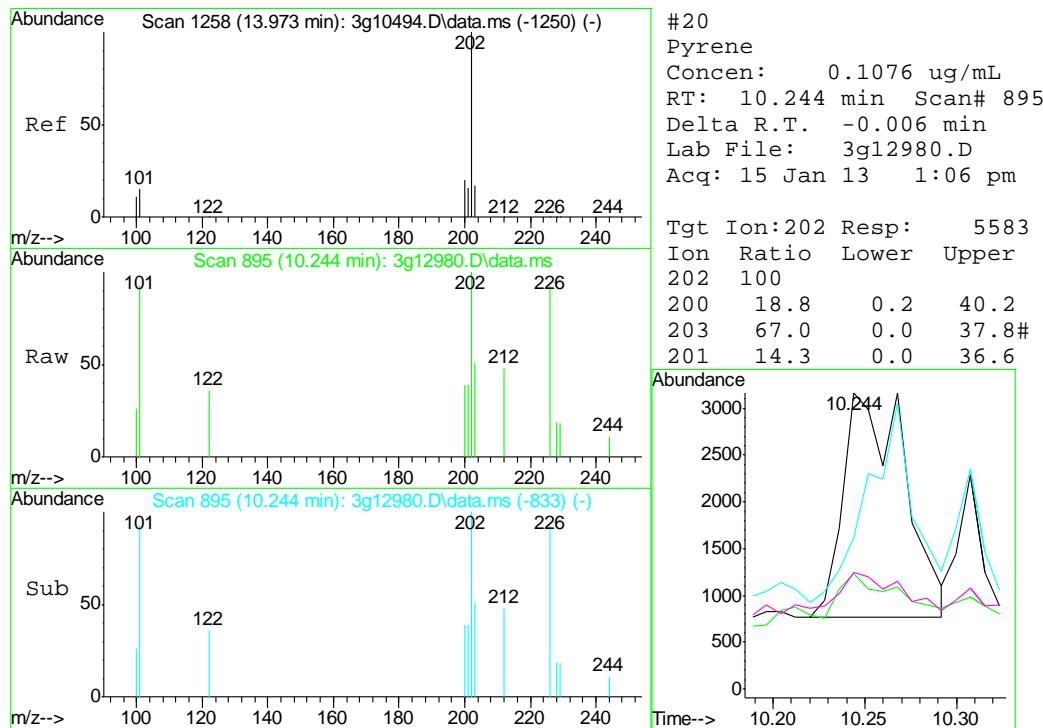
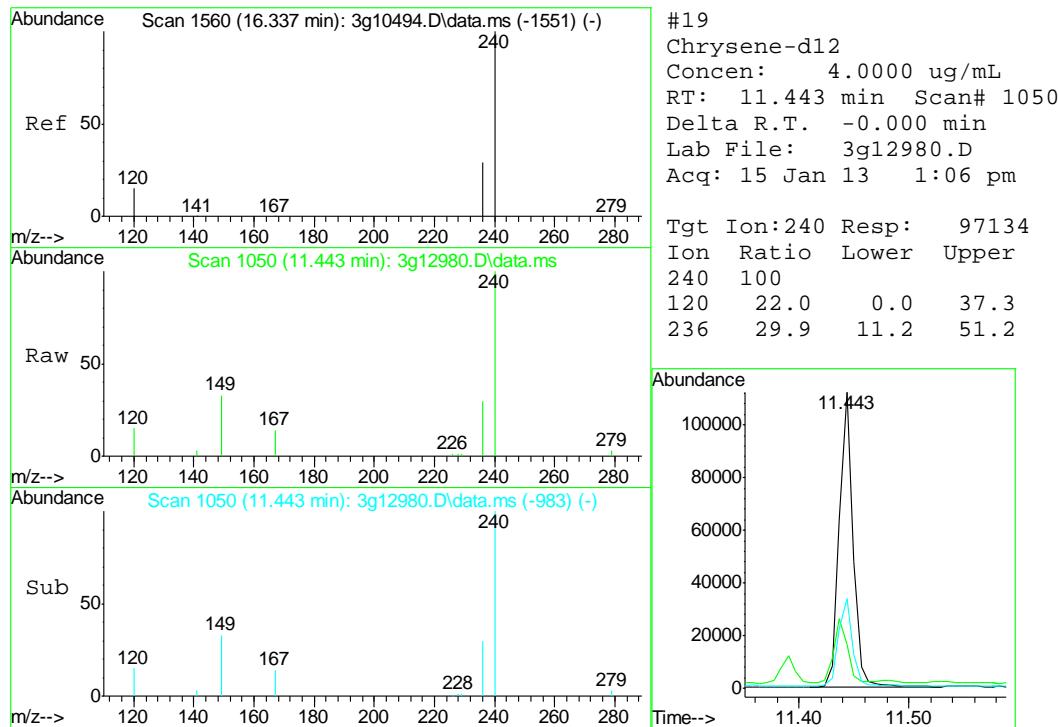


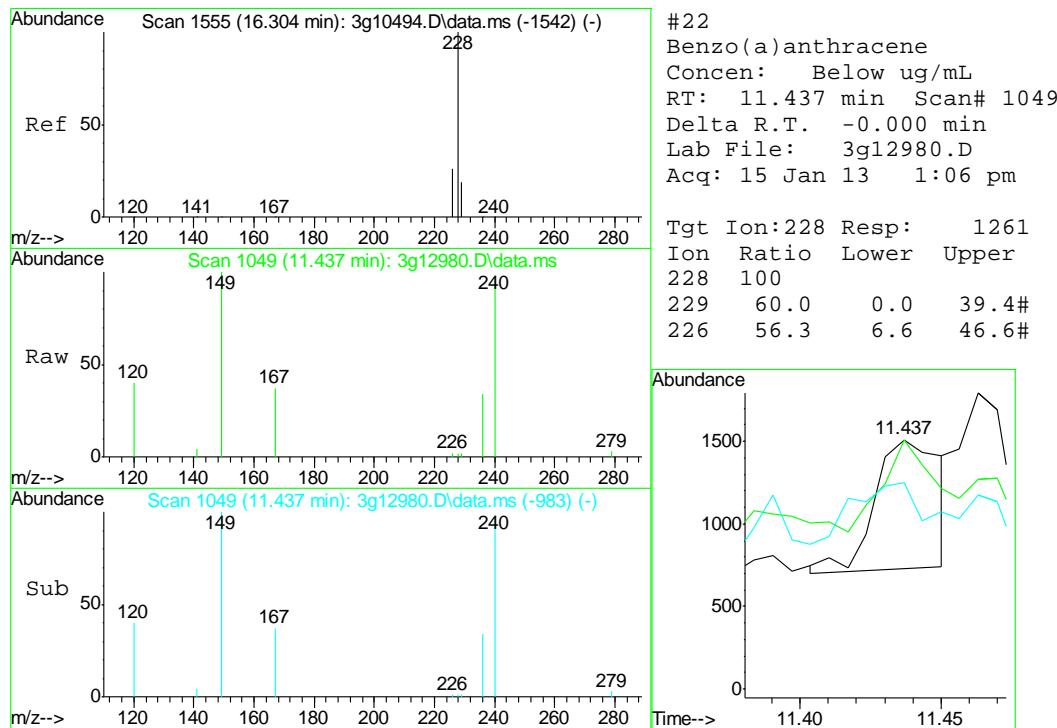
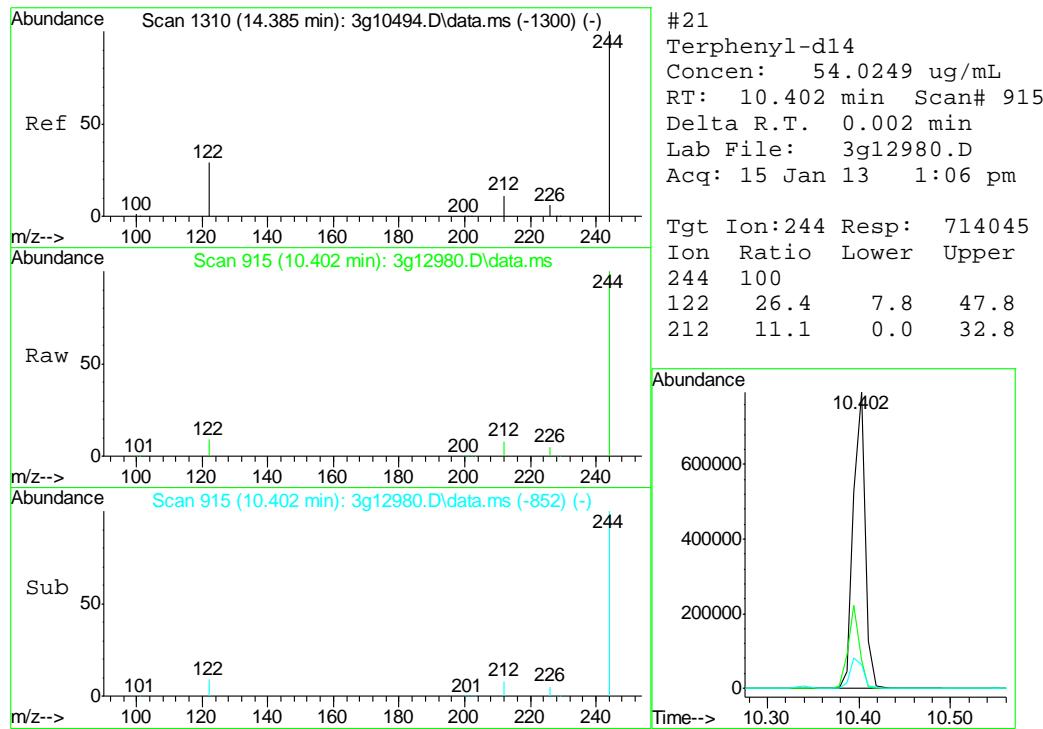


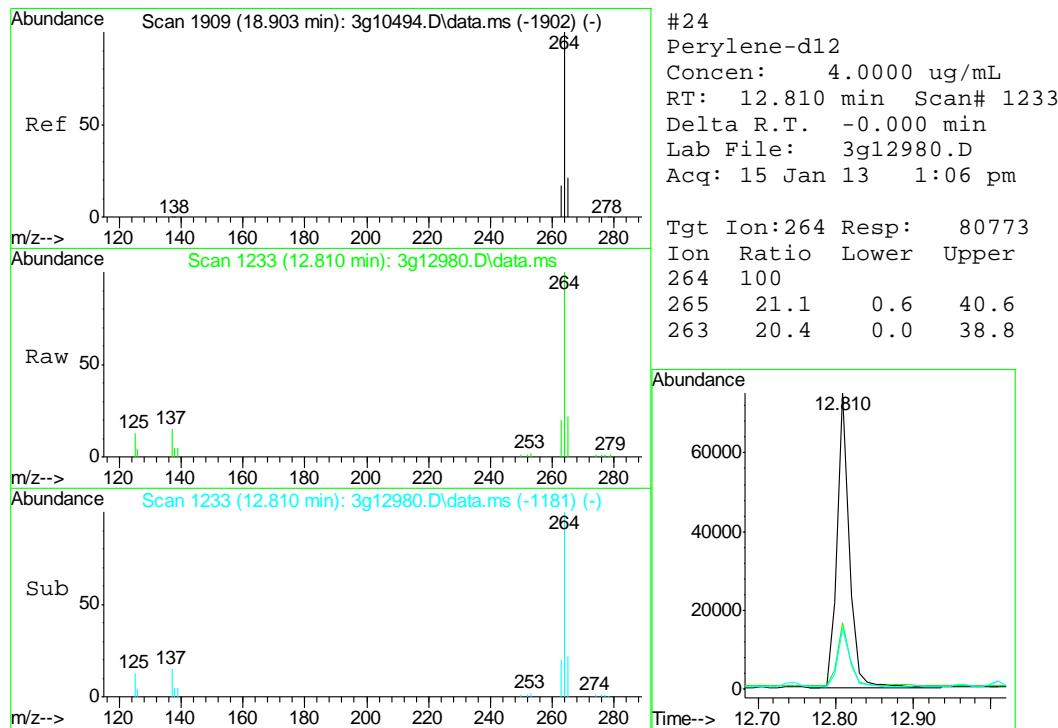
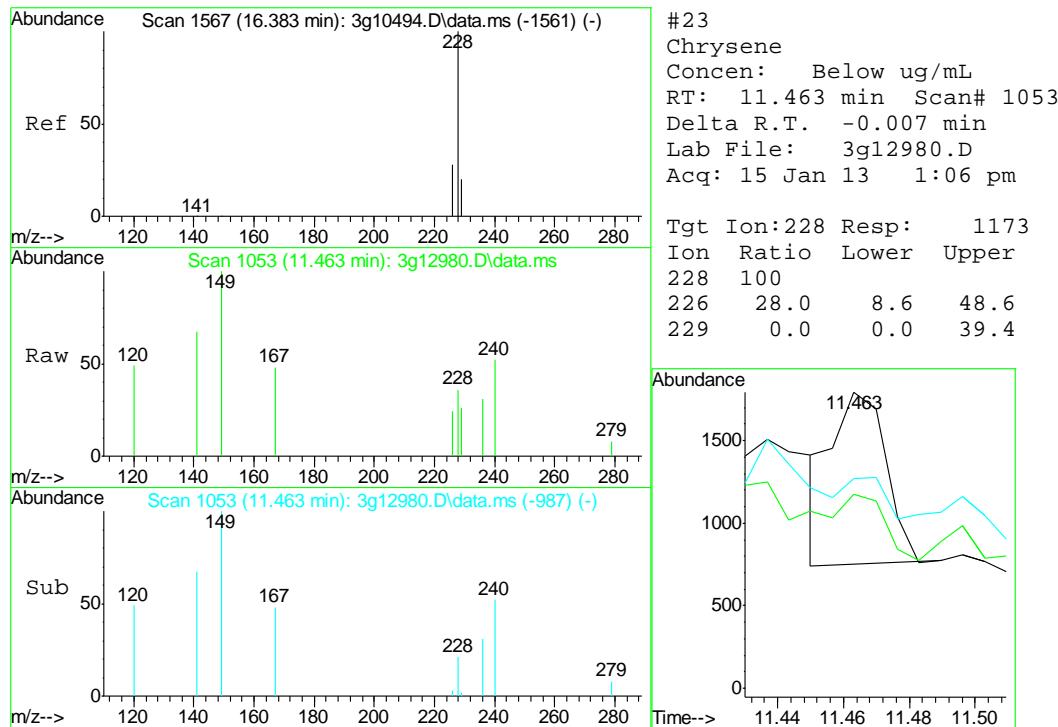


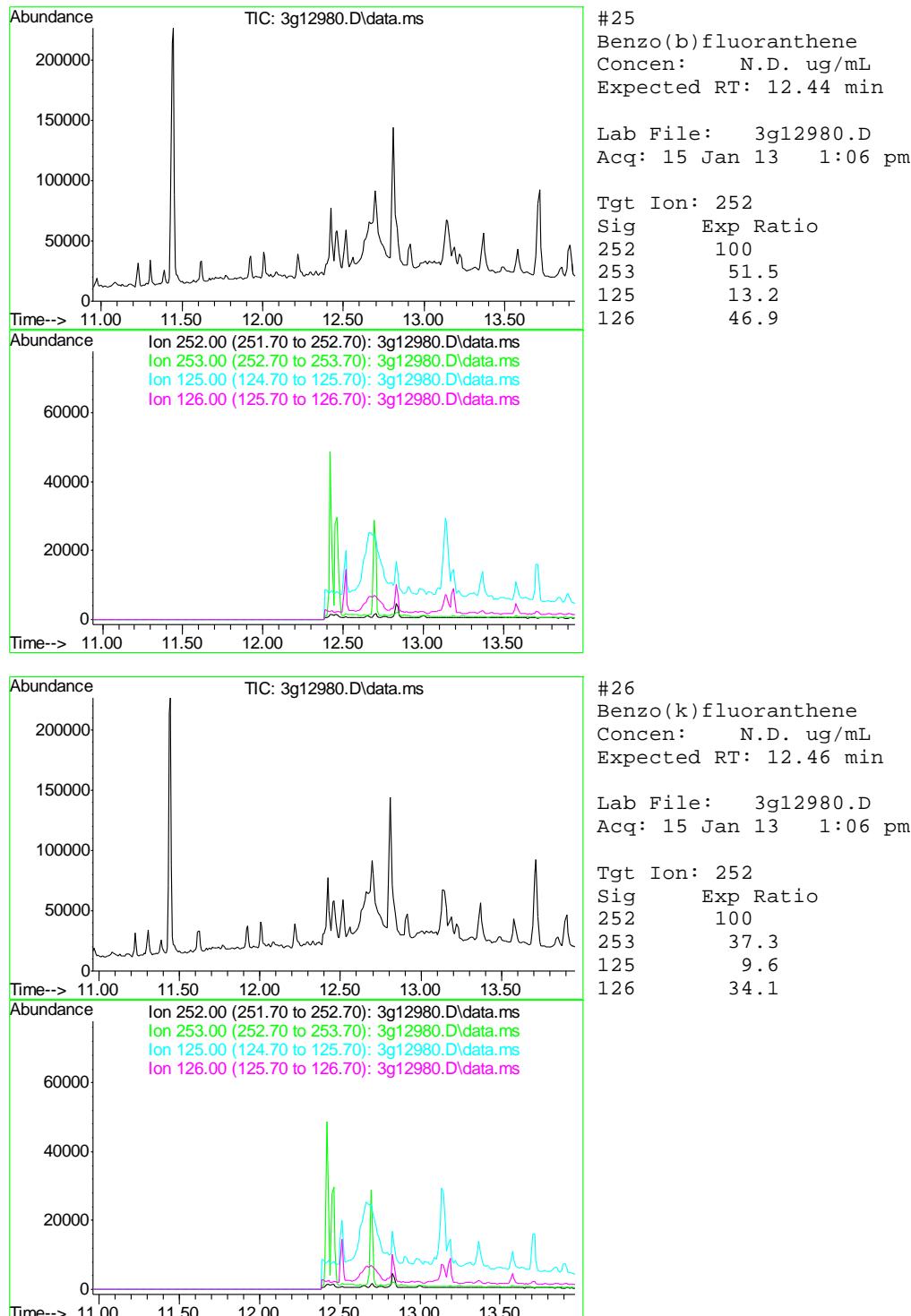


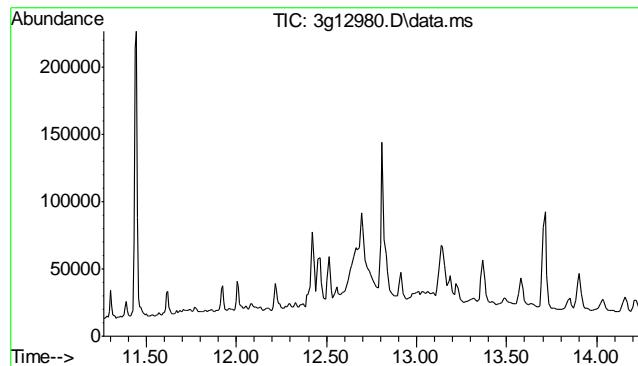








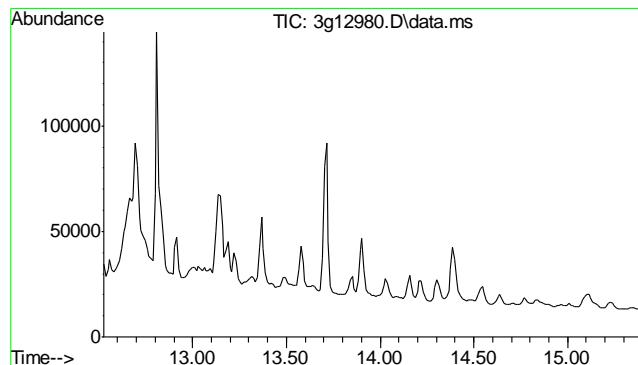
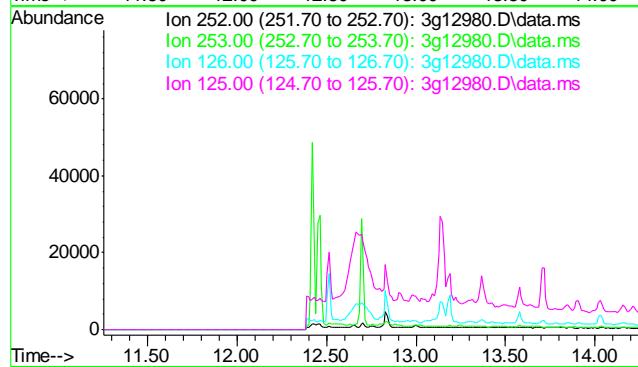




#27  
 Benzo(a)pyrene  
 Concen: N.D. ug/mL  
 Expected RT: 12.76 min

Lab File: 3g12980.D  
 Acq: 15 Jan 13 1:06 pm

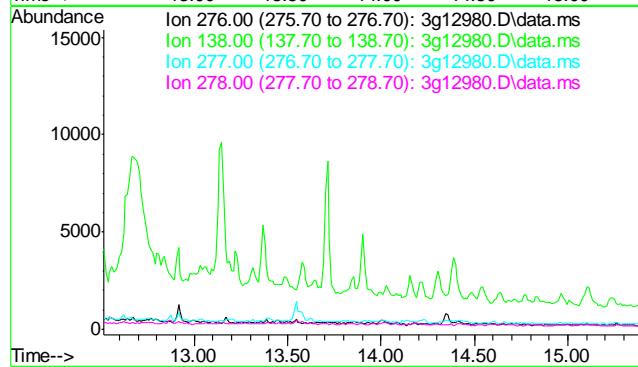
Tgt Ion:	Sig	Exp Ratio
252	100	
253	21.5	
126	20.4	
125	14.5	

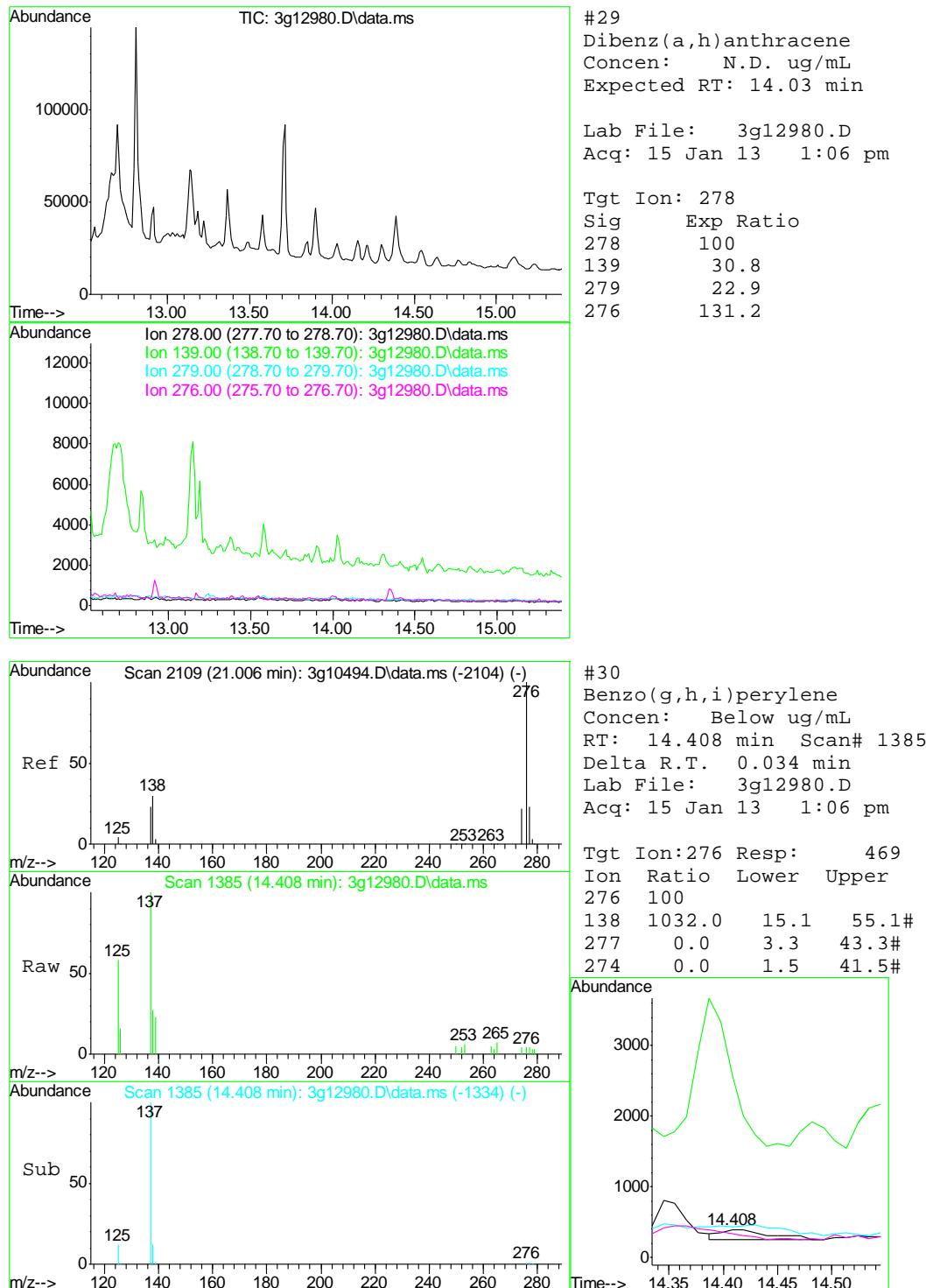


#28  
 Indeno(1,2,3-cd)pyrene  
 Concen: N.D. ug/mL  
 Expected RT: 14.02 min

Lab File: 3g12980.D  
 Acq: 15 Jan 13 1:06 pm

Tgt Ion:	Sig	Exp Ratio
276	100	
138	40.0	
277	24.8	
278	76.2	





## Quantitation Report (QT Reviewed)

Manual Integrations  
APPROVED  
(compounds with "m" flag)

Judy Nelson  
01/16/13 11:51

Data Path : C:\msdchem\1\DATA\011513\  
Data File : 3g12987.D  
Acq On : 15 Jan 2013 3:52 pm  
Operator : DONC  
Sample : D42556-2  
Misc : OP7223,E3G621,30.02,,,1,1  
ALS Vial : 18 Sample Multiplier: 1

Quant Time: Jan 16 09:11:53 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	5.620	136	162561	4.0000	ug/mL	0.00
6) Acenaphthene-d10	7.337	164	89868	4.0000	ug/mL	0.01
15) Phenanthrene-d10	8.819	188	154602	4.0000	ug/mL	0.00
19) Chrysene-d12	11.450	240	89179	4.0000	ug/mL	0.00
24) Perylene-d12	12.820	264	62316	4.0000	ug/mL	0.01

## System Monitoring Compounds

2) Nitrobenzene-d5	4.935	82	427246	29.2195	ug/mL	-0.01
Spiked Amount	50.000	Range	25 - 135	Recovery	=	58.44%
7) 2-Fluorobiphenyl	6.676	172	1417101	41.0897	ug/mL	0.00
Spiked Amount	50.000	Range	25 - 135	Recovery	=	82.18%
21) Terphenyl-d14	10.402	244	538734	44.3967	ug/mL	0.00
Spiked Amount	50.000	Range	25 - 135	Recovery	=	88.80%

## Target Compounds

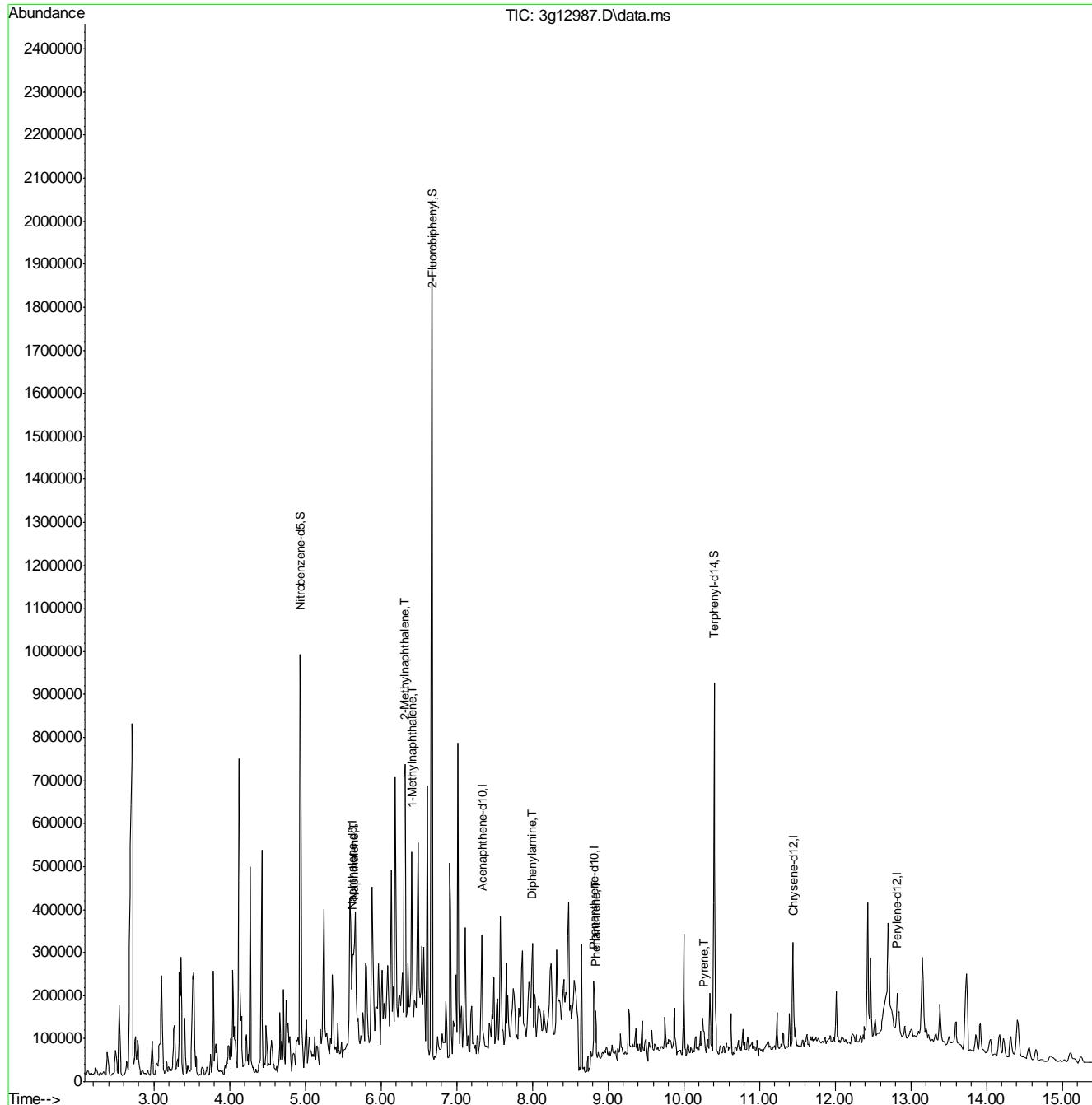
				Qvalue
3) N-Nitrosodimethylamine	0.000	74	0	N.D. d
4) N-Nitrosodi-propylamine	0.000	70	0	N.D. d
5) Naphthalene	5.645	128	212128	4.5354 ug/mL 72
8) 2-Methylnaphthalene	6.319	142	378787	13.1167 ug/mL 93
9) 1-Methylnaphthalene	6.406	142	155245m	6.1464 ug/mL
10) Acenaphthylene	0.000	152	0	N.D. d
11) Acenaphthene	0.000	154	0	N.D. d
12) Dibenzofuran	0.000	168	0	N.D. d
13) Fluorene	0.000	166	0	N.D. d
14) Diphenylamine	7.999	169	142230m	5.0440 ug/mL
16) Phenanthrene	8.835	178	75722	1.2674 ug/mL# 44
17) Anthracene	0.000	178	0	N.D. d
18) Fluoranthene	0.000	202	0	N.D. d
20) Pyrene	10.268	202	31604	0.6636 ug/mL# 58
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	0.000	276	0	N.D. d

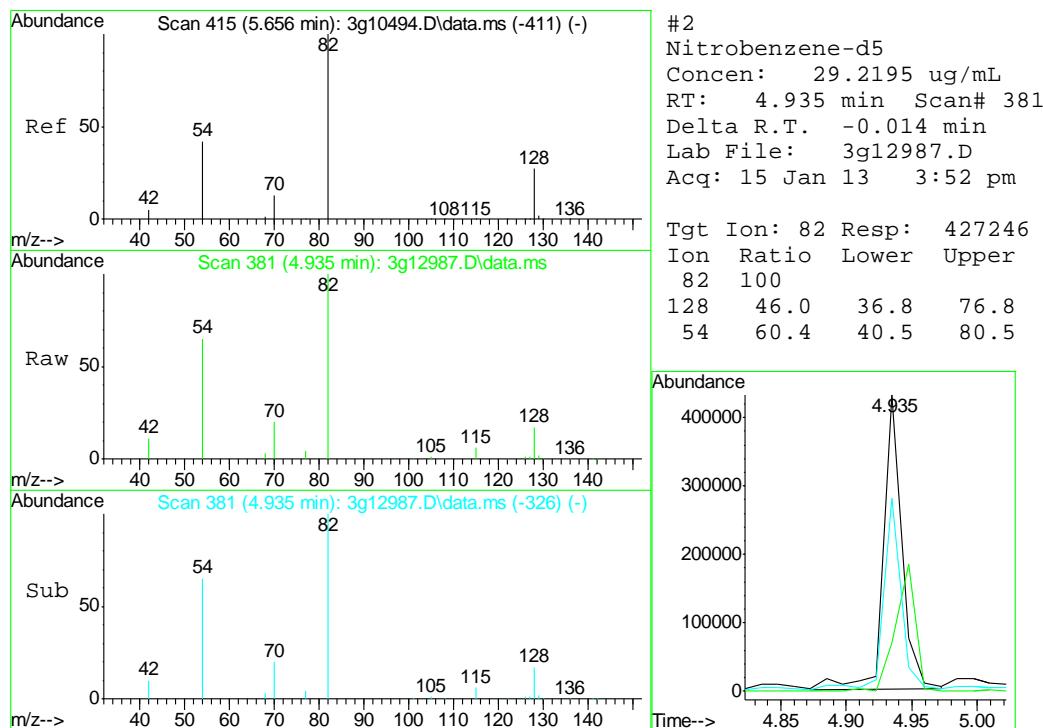
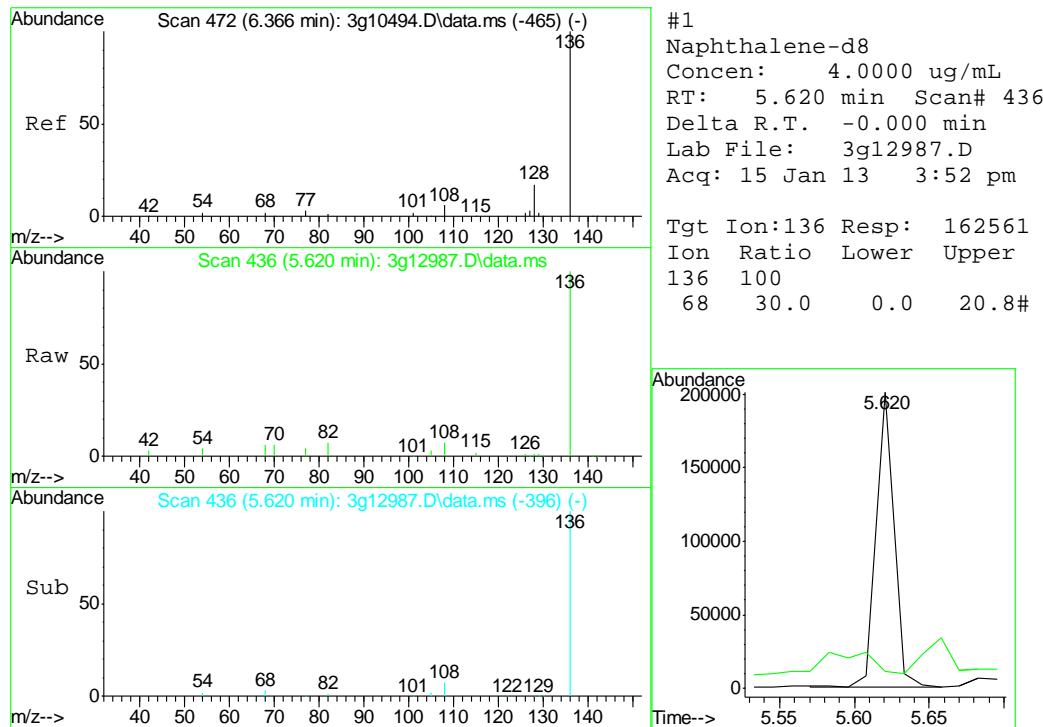
(#) = qualifier out of range (m) = manual integration (+) = signals summed

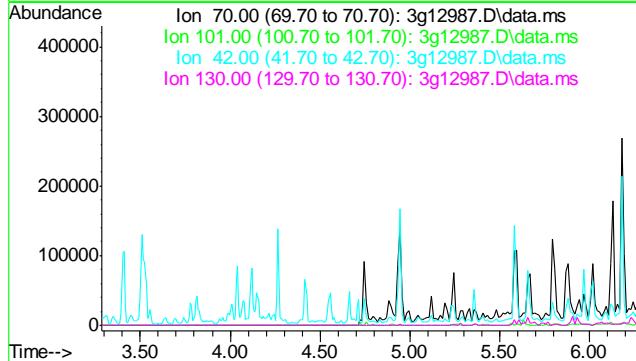
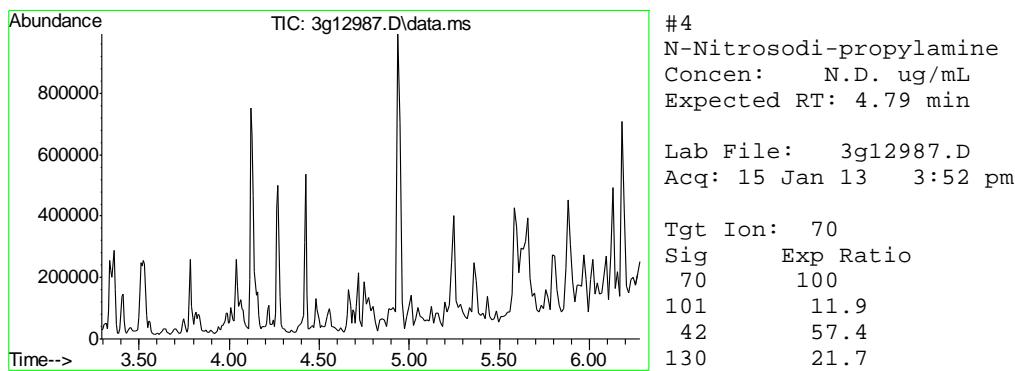
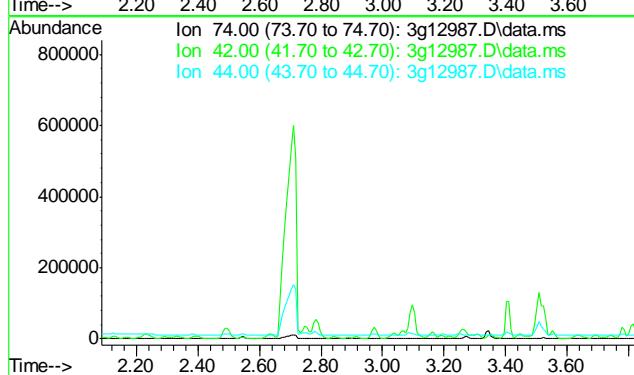
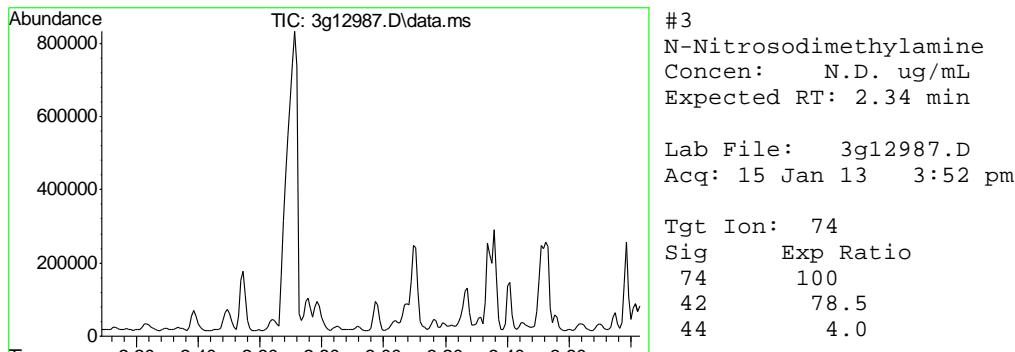
## Quantitation Report (QT Reviewed)

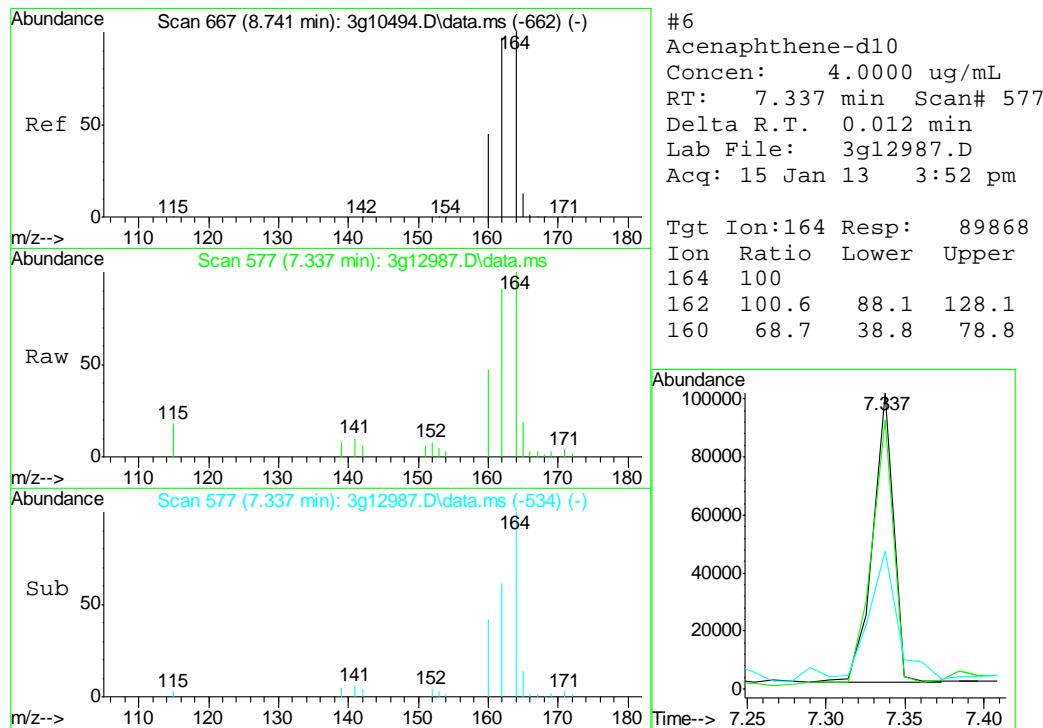
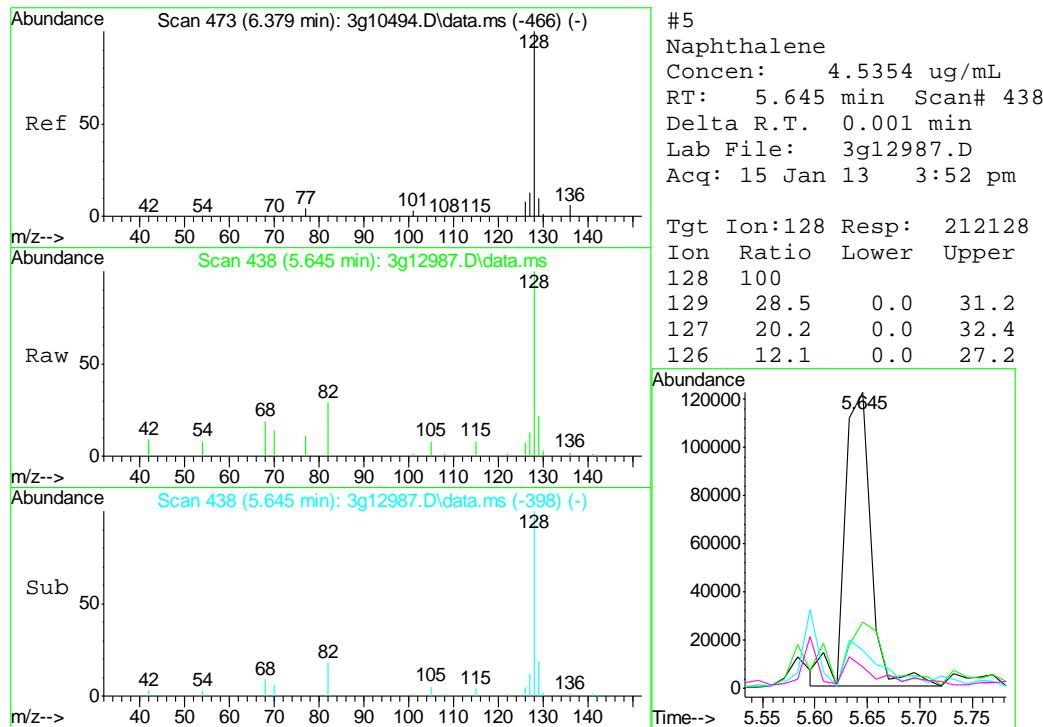
Data Path : C:\msdchem\1\DATA\011513\  
 Data File : 3g12987.D  
 Acq On : 15 Jan 2013 3:52 pm  
 Operator : DONC  
 Sample : D42556-2  
 Misc : OP7223,E3G621,30.02,,,1,1  
 ALS Vial : 18 Sample Multiplier: 1

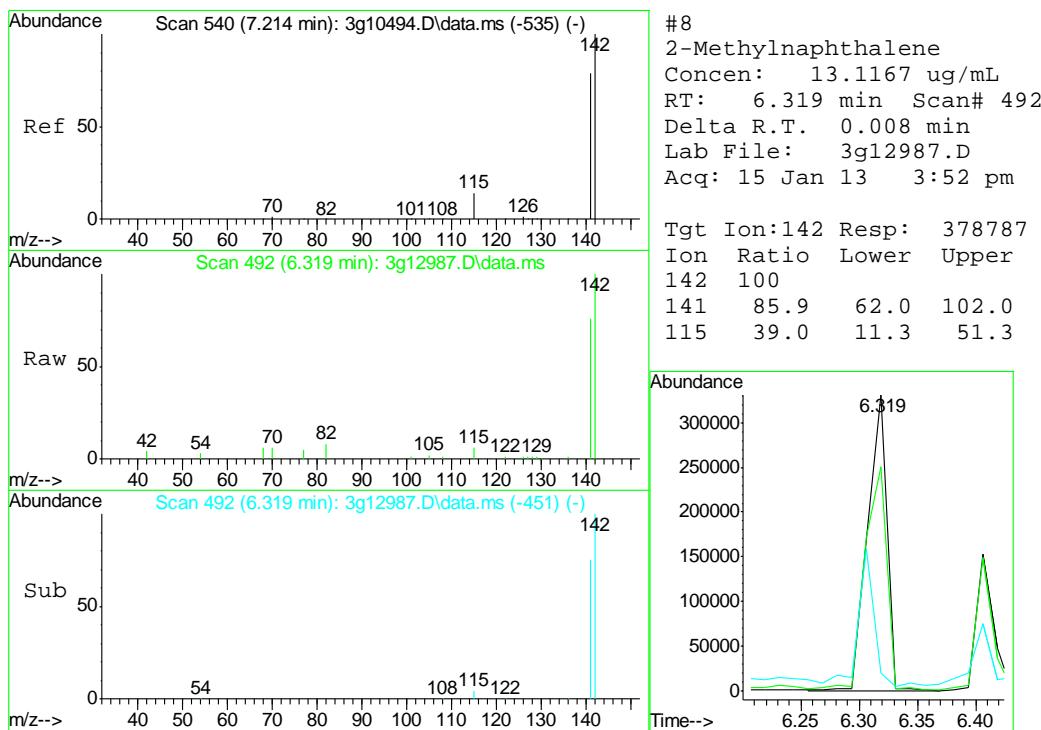
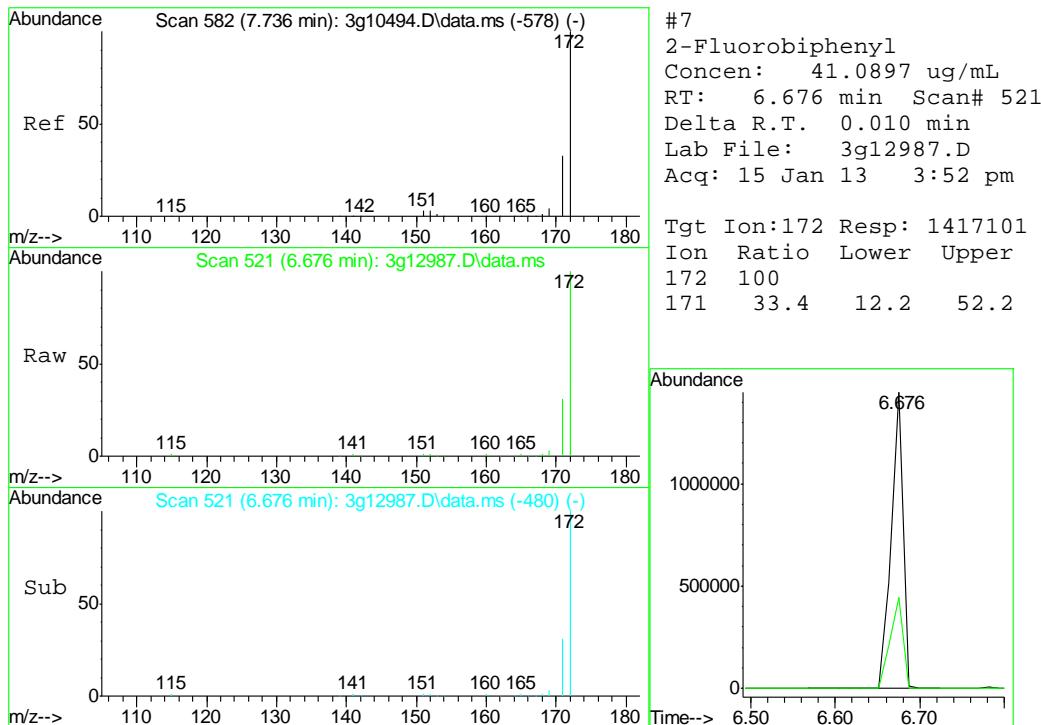
Quant Time: Jan 16 09:11:53 2013  
 Quant Method : C:\msdchem\1\METHODS\SIMPE3G611.M  
 Quant Title : PAHSIM BASE  
 QLast Update : Thu Jan 10 14:18:35 2013  
 Response via : Initial Calibration

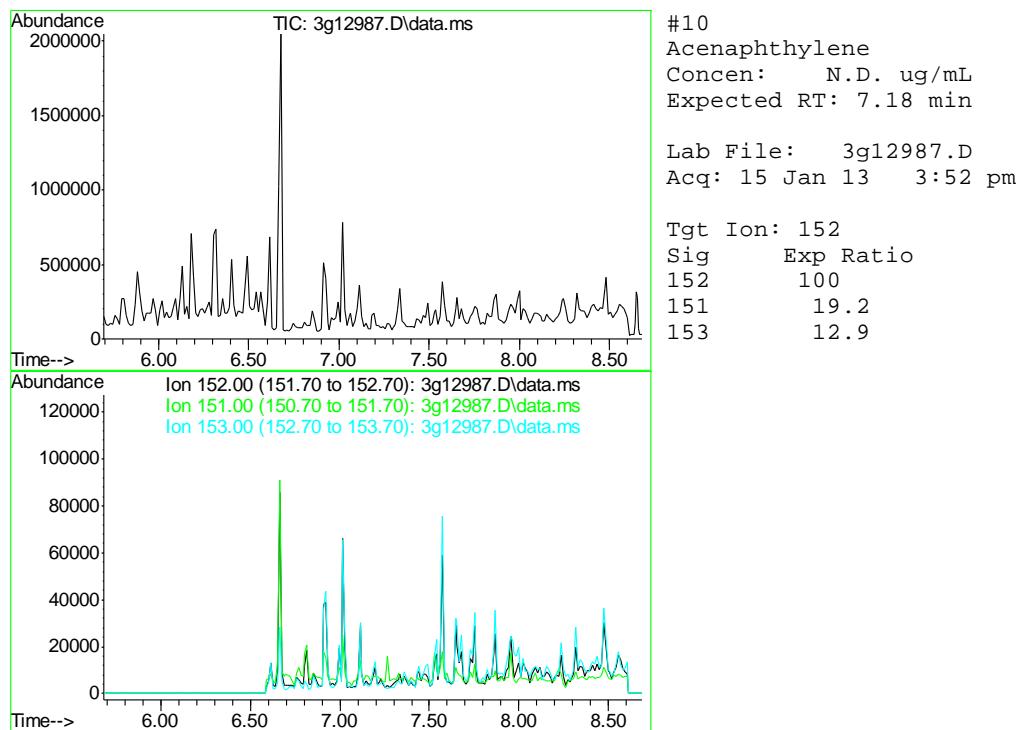
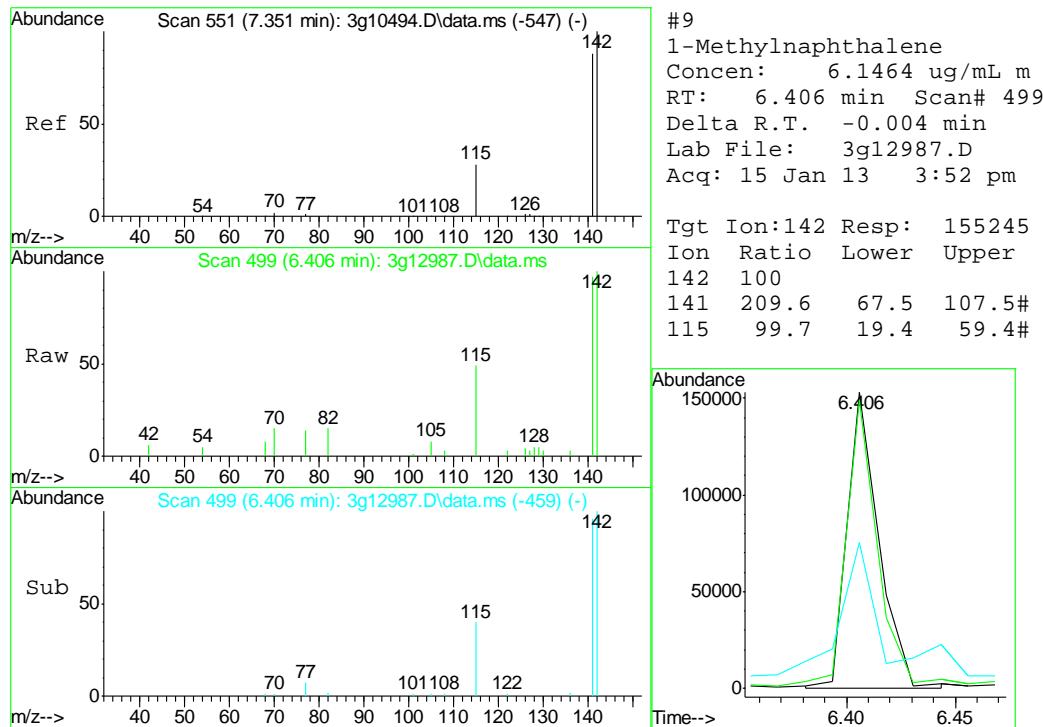


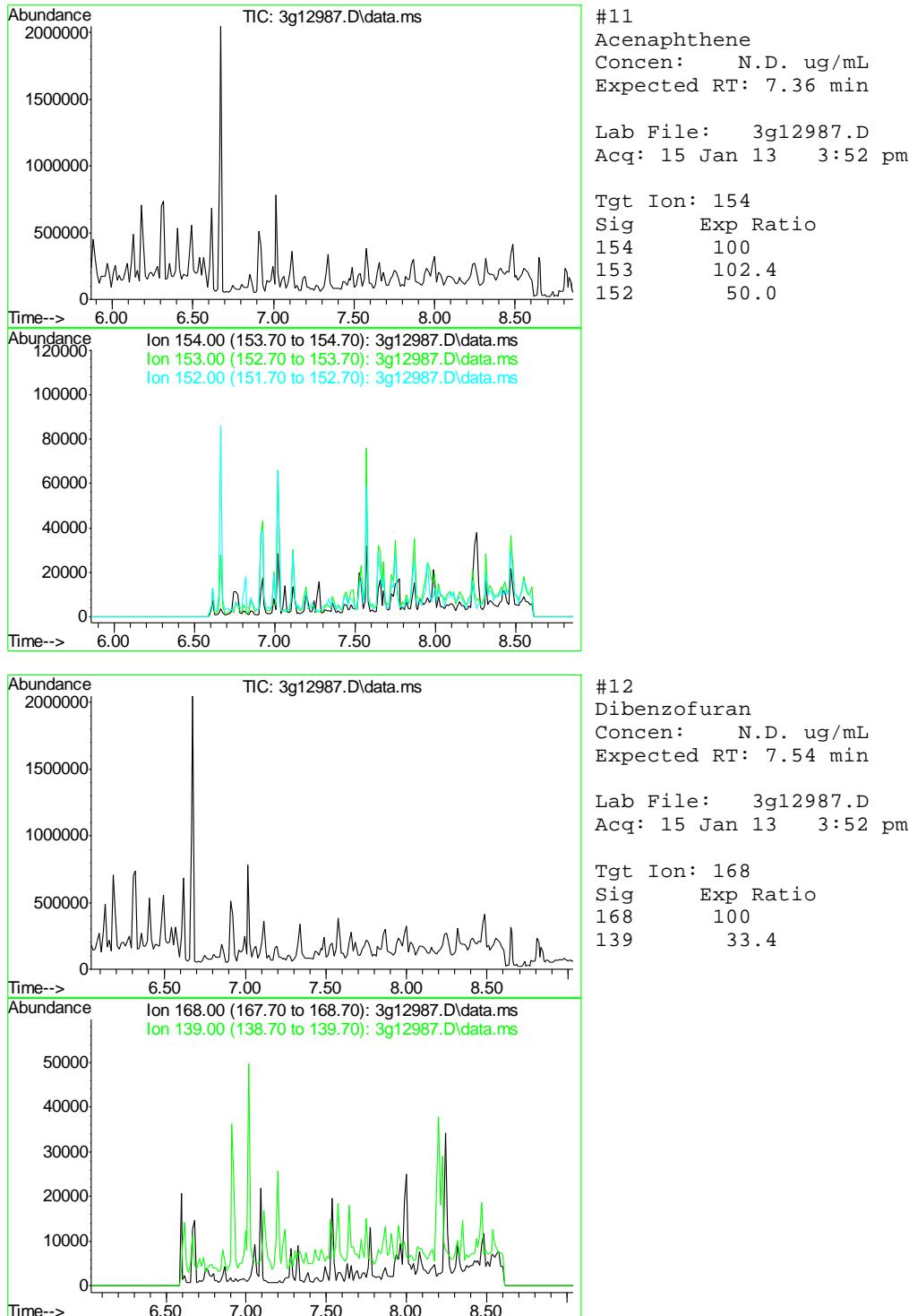


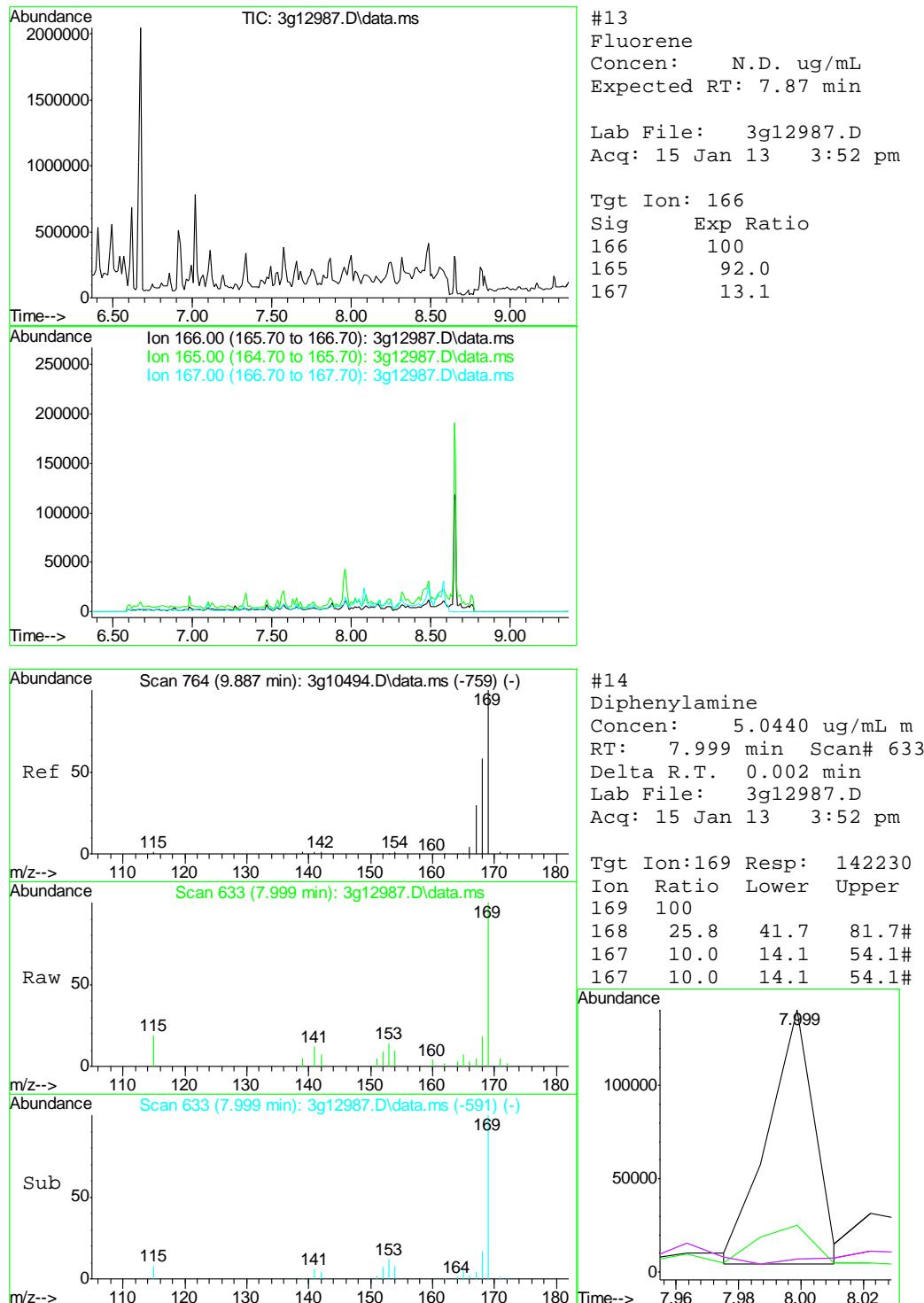


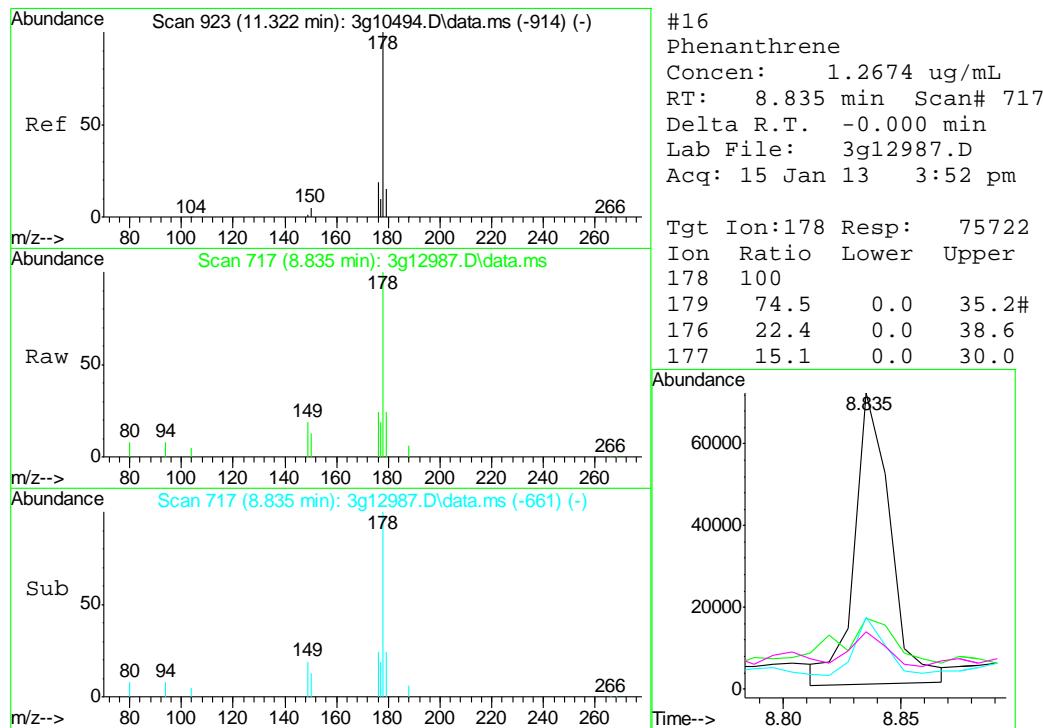
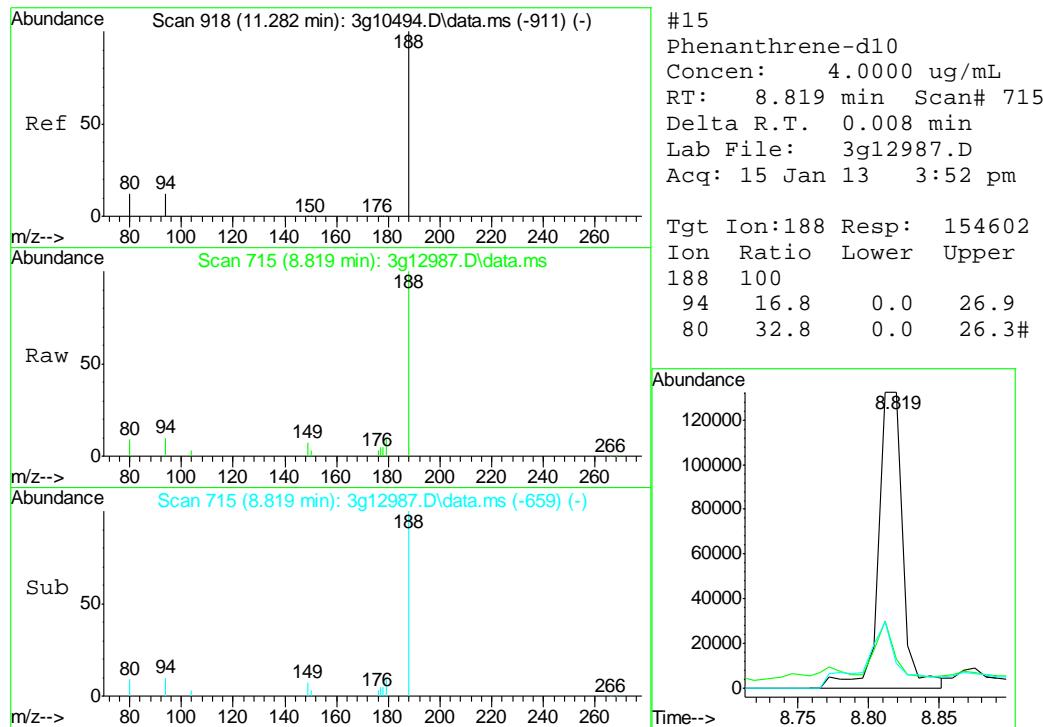


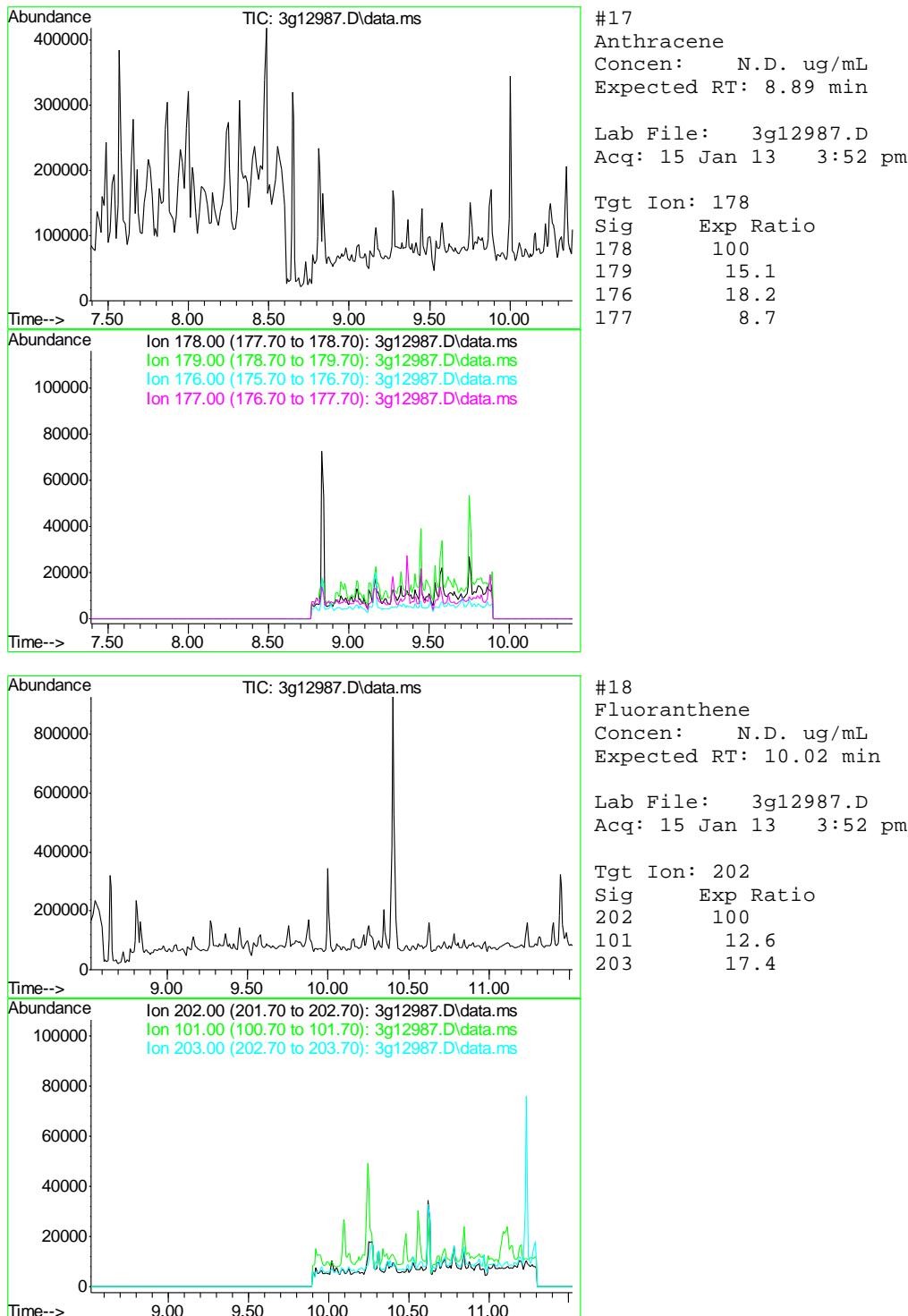


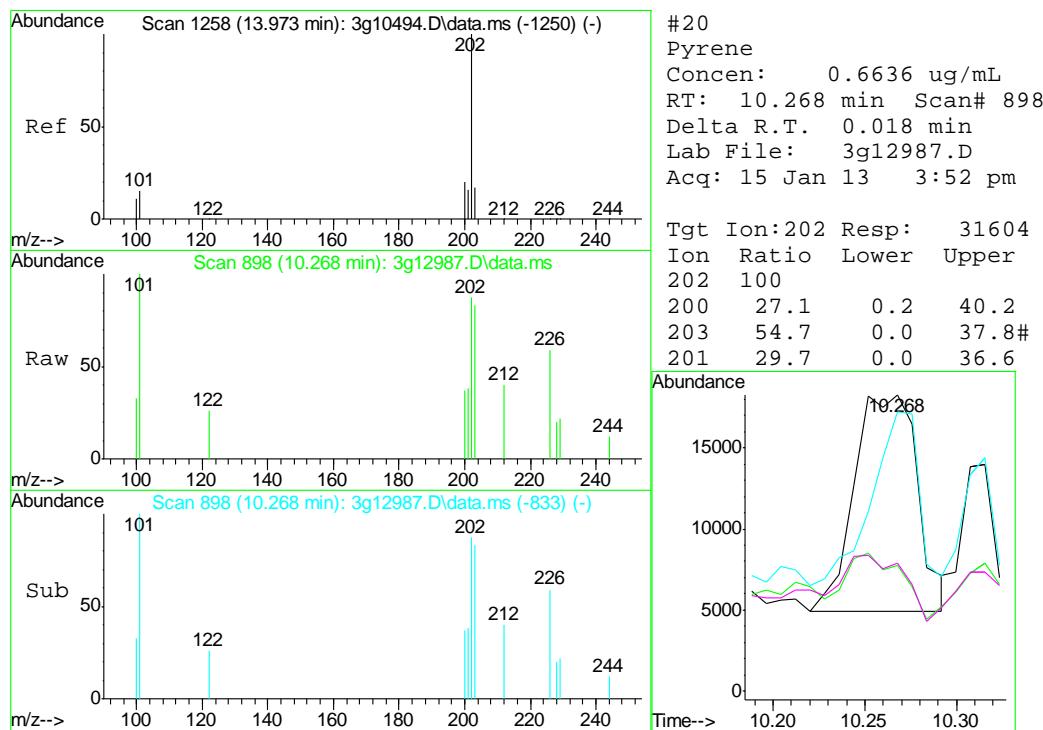
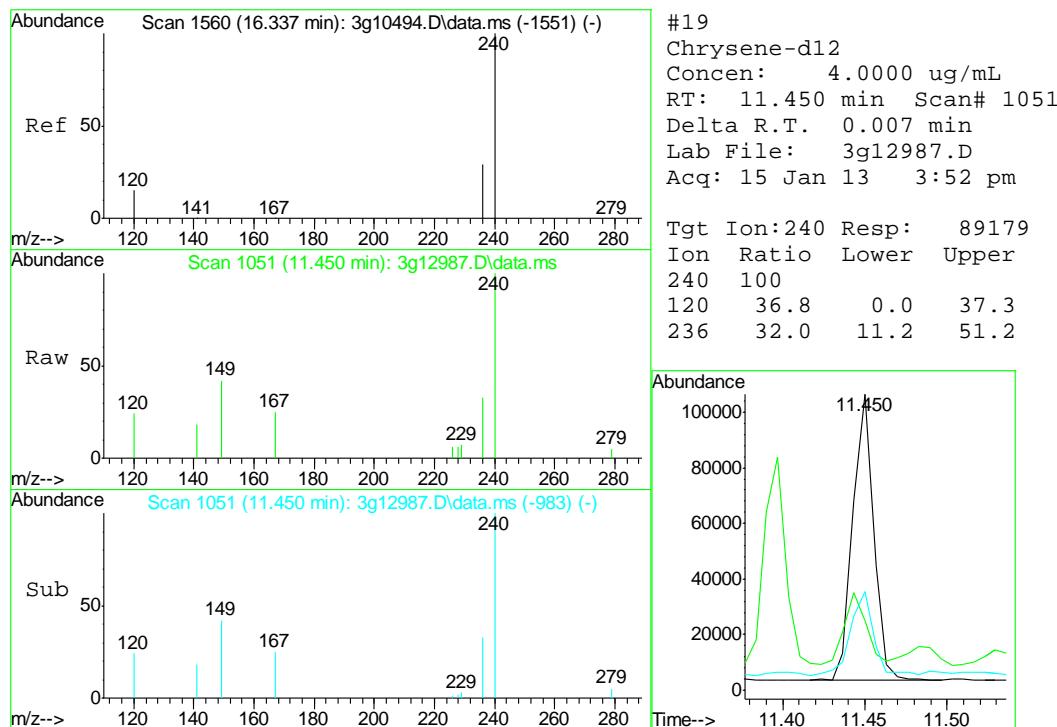


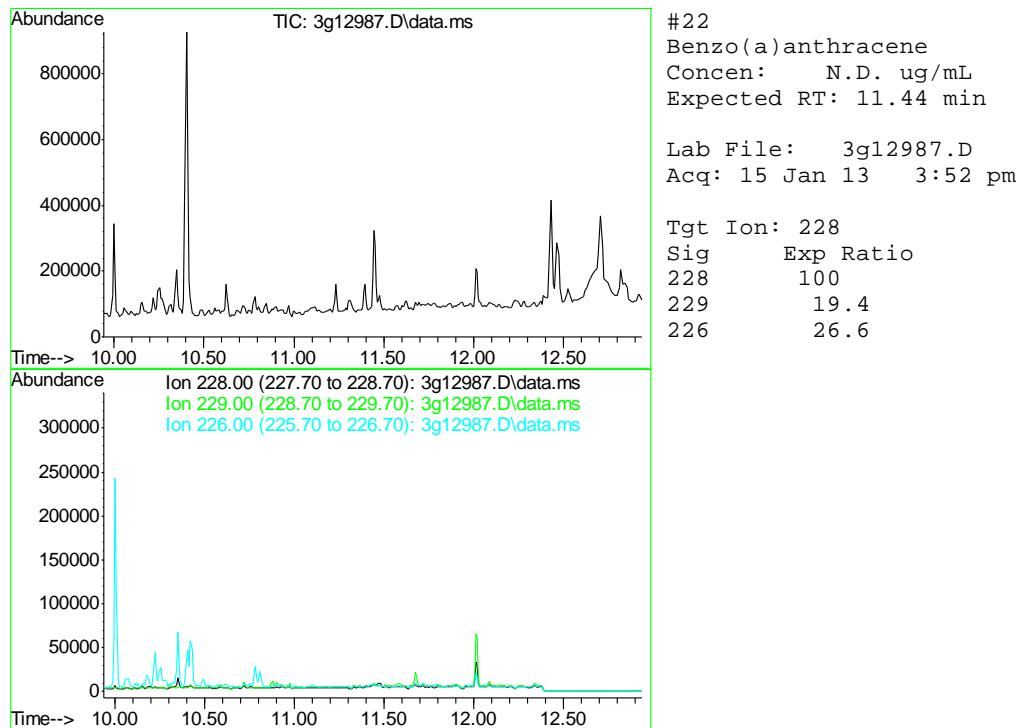
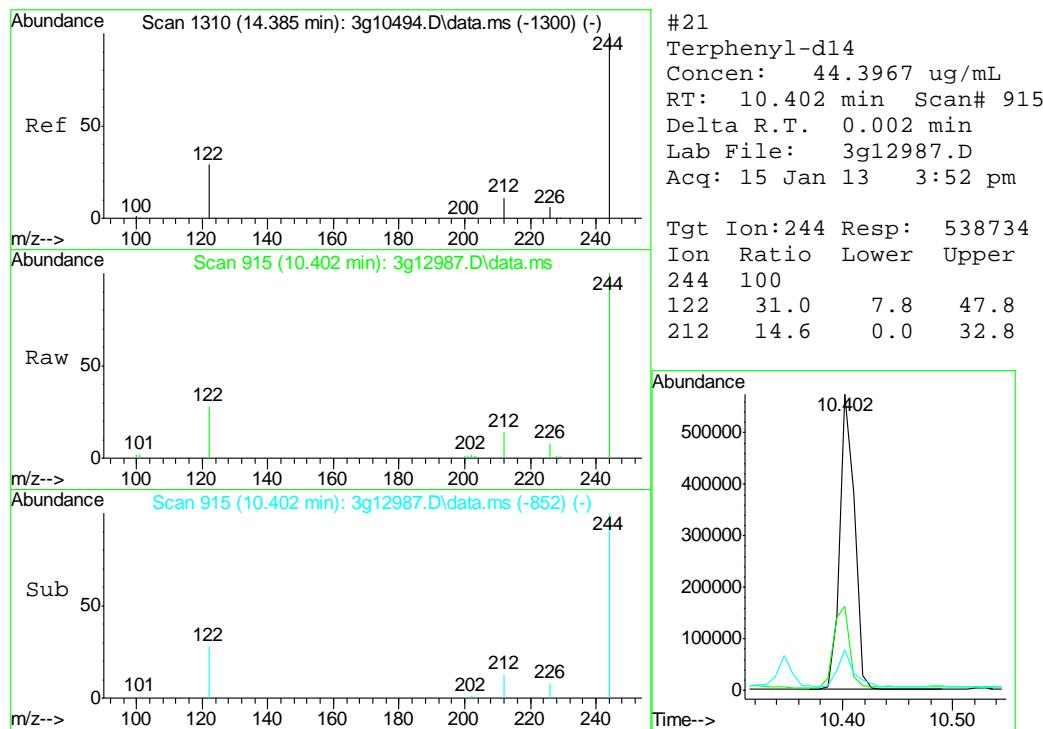


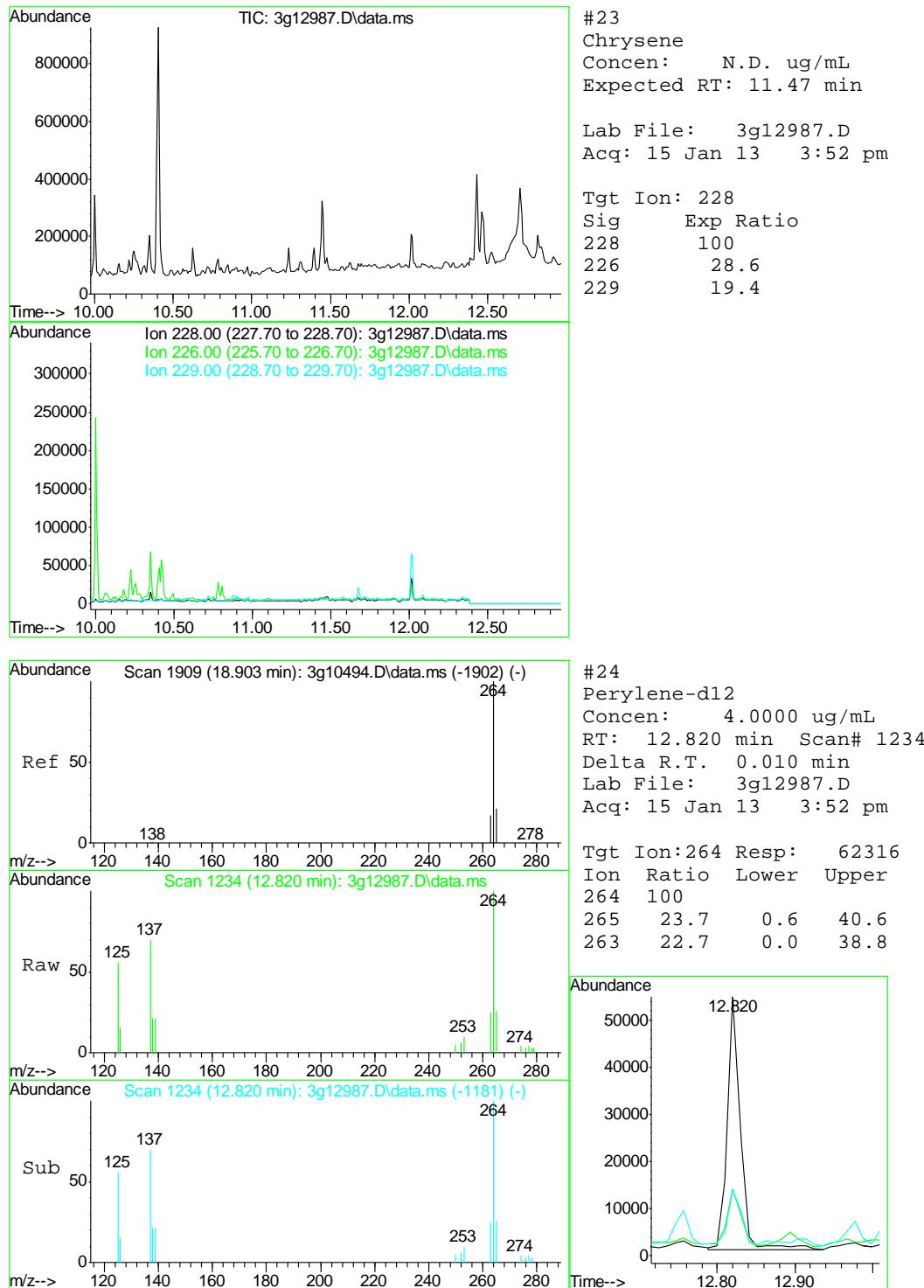


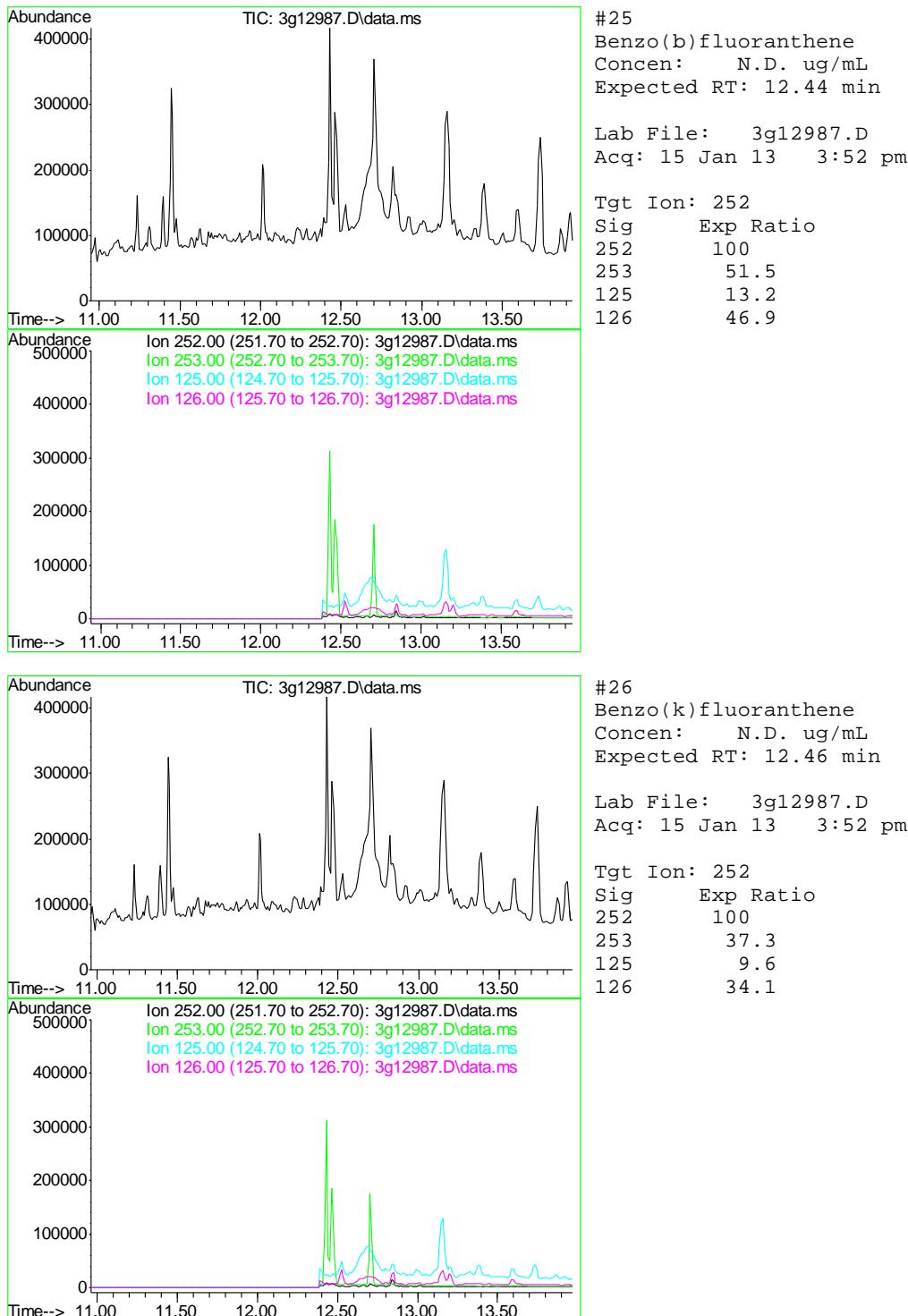


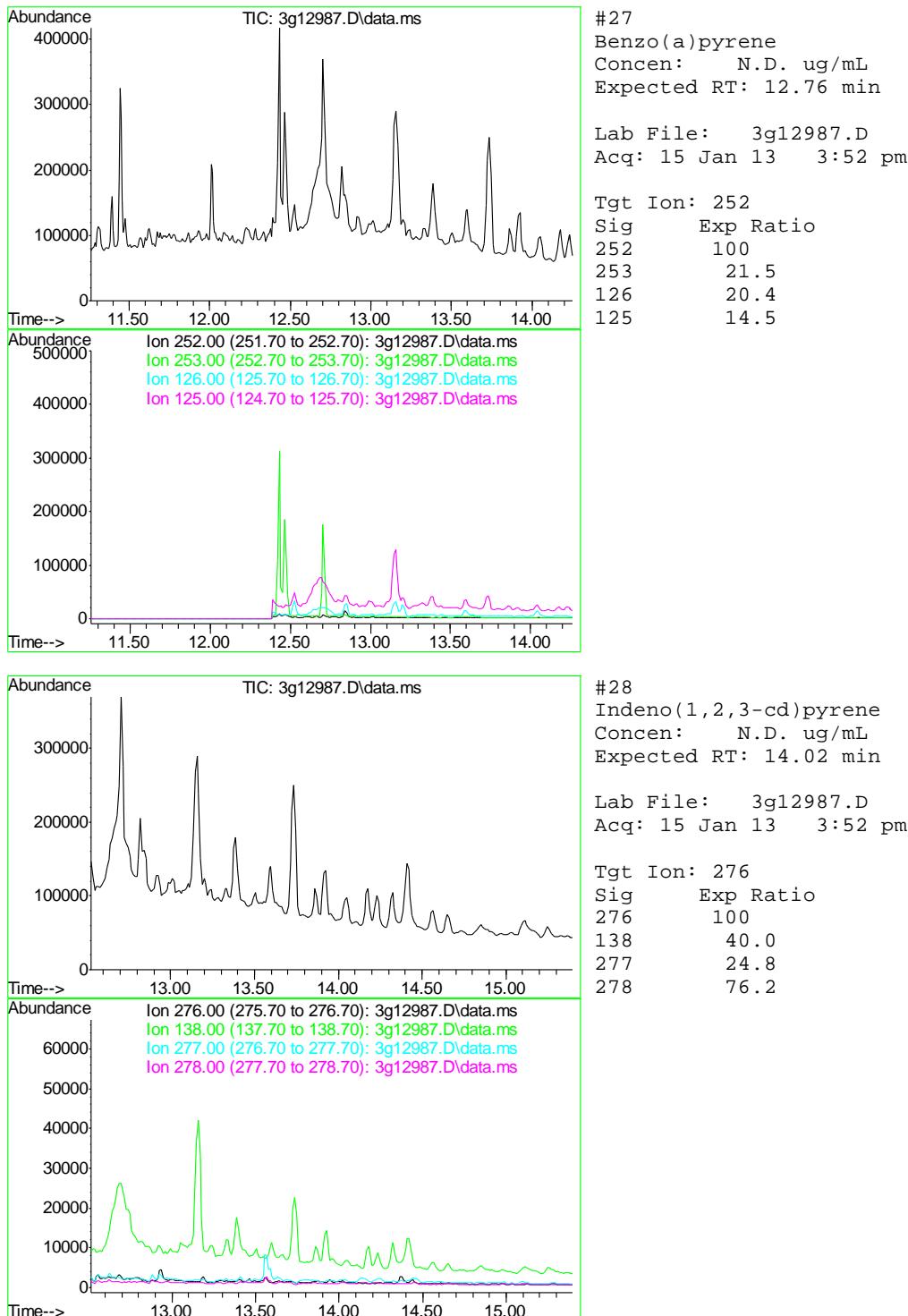


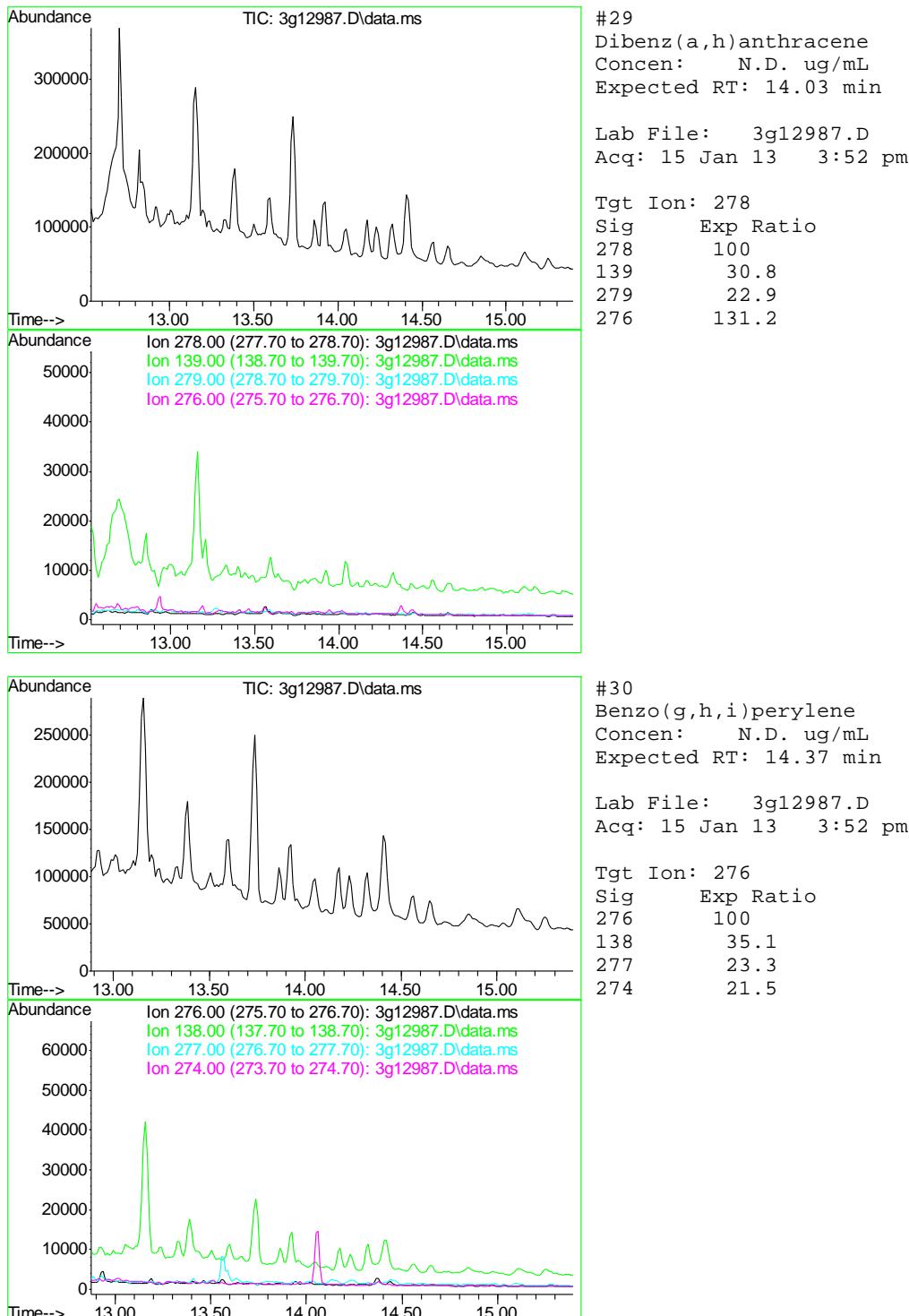












Judy Nelson  
 01/16/13 11:51

## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\011513\  
 Data File : 3g12973.D  
 Acq On : 15 Jan 2013 10:14 am  
 Operator : DONC  
 Sample : OP7223-MB  
 Misc : OP7223,E3G621,30.00,,,1,1  
 ALS Vial : 4 Sample Multiplier: 1

Quant Time: Jan 15 13:46:44 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	122903	4.0000	ug/mL	0.00
6) Acenaphthene-d10	7.337	164	71901	4.0000	ug/mL	0.01
15) Phenanthrene-d10	8.812	188	128999	4.0000	ug/mL	0.00
19) Chrysene-d12	11.443	240	101544	4.0000	ug/mL	0.00
24) Perylene-d12	12.810	264	84293	4.0000	ug/mL	0.00

## System Monitoring Compounds

2) Nitrobenzene-d5	4.935	82	445750	40.3218	ug/mL	-0.01
Spiked Amount	50.000	Range	25 - 135	Recovery	=	80.64%
7) 2-Fluorobiphenyl	6.676	172	1170373	42.6075	ug/mL	0.00
Spiked Amount	50.000	Range	25 - 135	Recovery	=	85.22%
21) Terphenyl-d14	10.402	244	725496	52.5073	ug/mL	0.00
Spiked Amount	50.000	Range	25 - 135	Recovery	=	105.02%

## Target Compounds

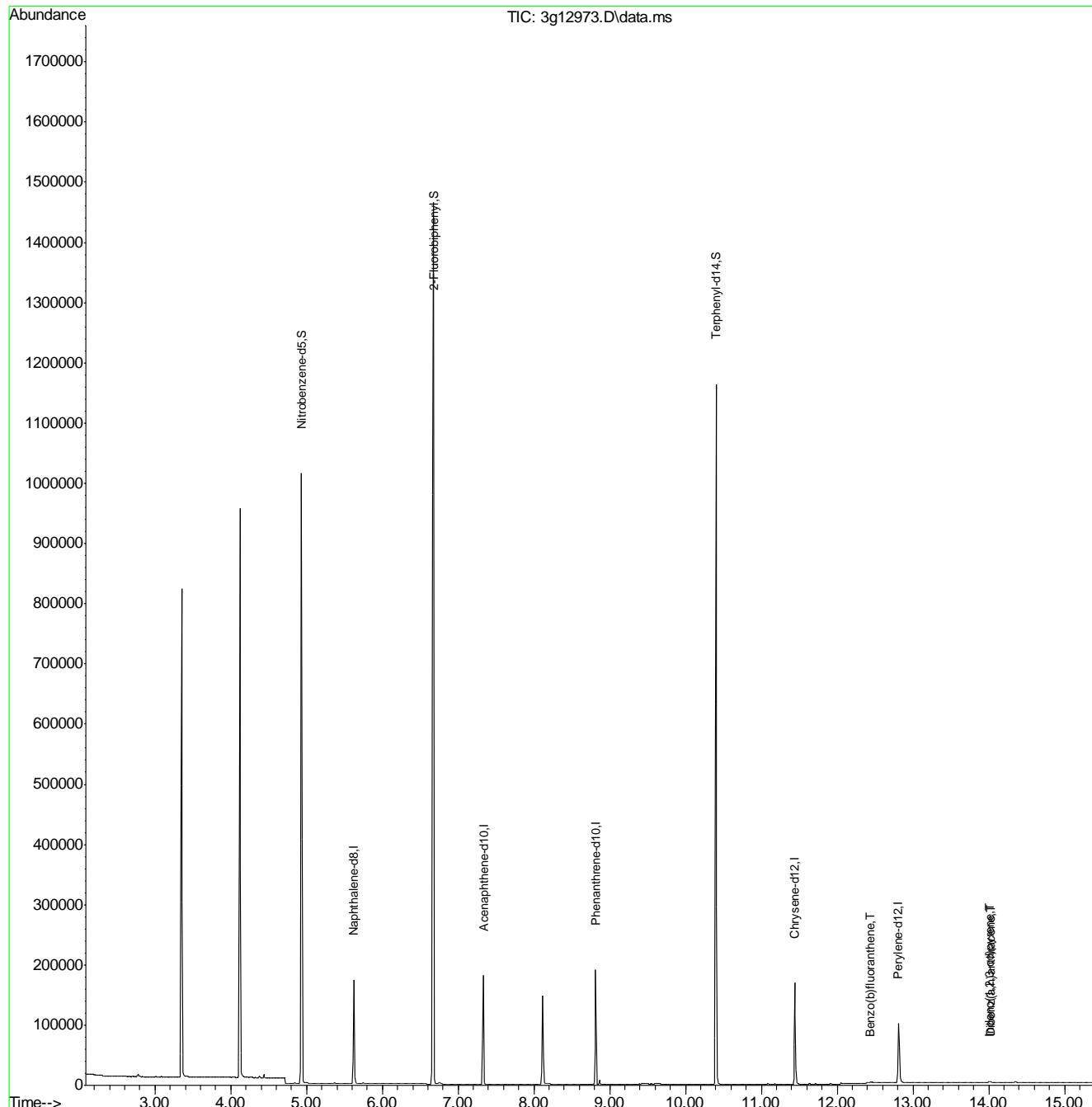
				Qvalue
3) N-Nitrosodimethylamine	2.356	74	24	N.D.
4) N-Nitrosodi-propylamine	0.000	70	0	N.D. d
5) Naphthalene	5.646	128	183	Below Cal # 48
8) 2-Methylnaphthalene	6.319	142	186	N.D.
9) 1-Methylnaphthalene	6.394	142	78	N.D.
10) Acenaphthylene	7.396	152	202	N.D.
11) Acenaphthene	7.385	154	61	N.D.
12) Dibenzofuran	7.857	168	73	N.D.
13) Fluorene	0.000	166	0	N.D. d
14) Diphenylamine	0.000	169	0	N.D. d
16) Phenanthrene	8.812	178	152	N.D.
17) Anthracene	8.891	178	39	N.D.
18) Fluoranthene	10.015	202	205	N.D.
20) Pyrene	10.244	202	152	N.D.
22) Benzo(a)anthracene	11.437	228	1041	N.D.
23) Chrysene	11.470	228	877	N.D.
25) Benzo(b)fluoranthene	12.442	252	1135m	0.0649 ug/mL
26) Benzo(k)fluoranthene	12.463	252	1206	N.D.
27) Benzo(a)pyrene	0.000	252	0	N.D. d
28) Indeno(1,2,3-cd)pyrene	14.009	276	1461	0.0777 ug/mL 89
29) Dibenz(a,h)anthracene	14.019	278	1174	0.0897 ug/mL 92
30) Benzo(g,h,i)perylene	14.356	276	1418	N.D.

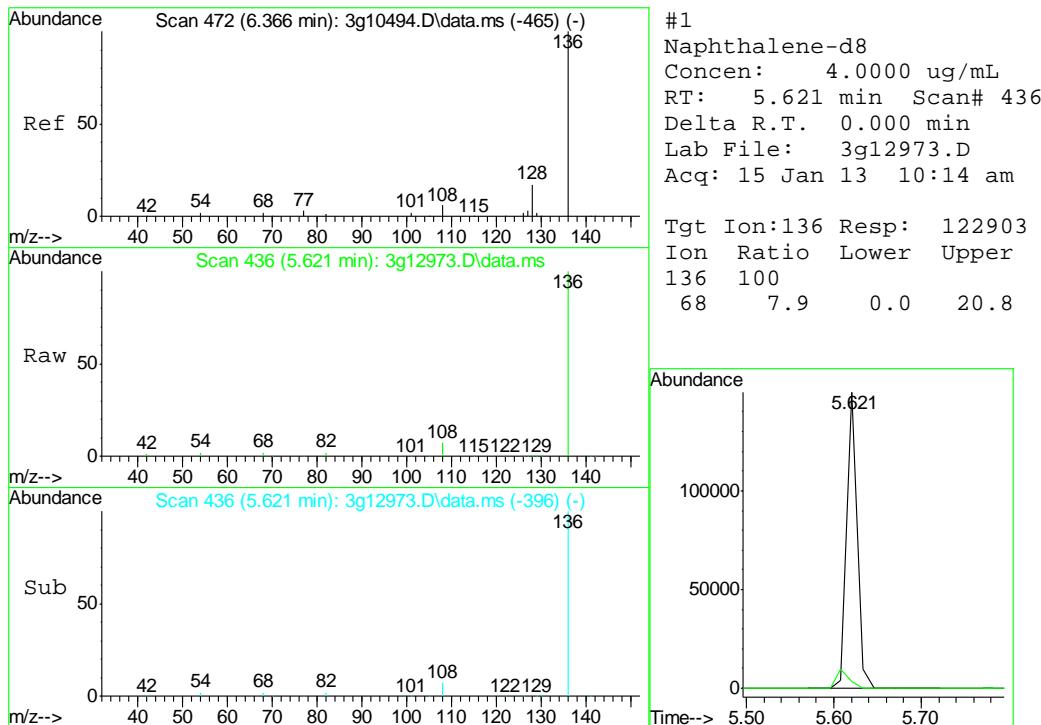
(#) = qualifier out of range (m) = manual integration (+) = signals summed

## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\011513\  
 Data File : 3g12973.D  
 Acq On : 15 Jan 2013 10:14 am  
 Operator : DONC  
 Sample : OP7223-MB  
 Misc : OP7223,E3G621,30.00,,,1,1  
 ALS Vial : 4 Sample Multiplier: 1

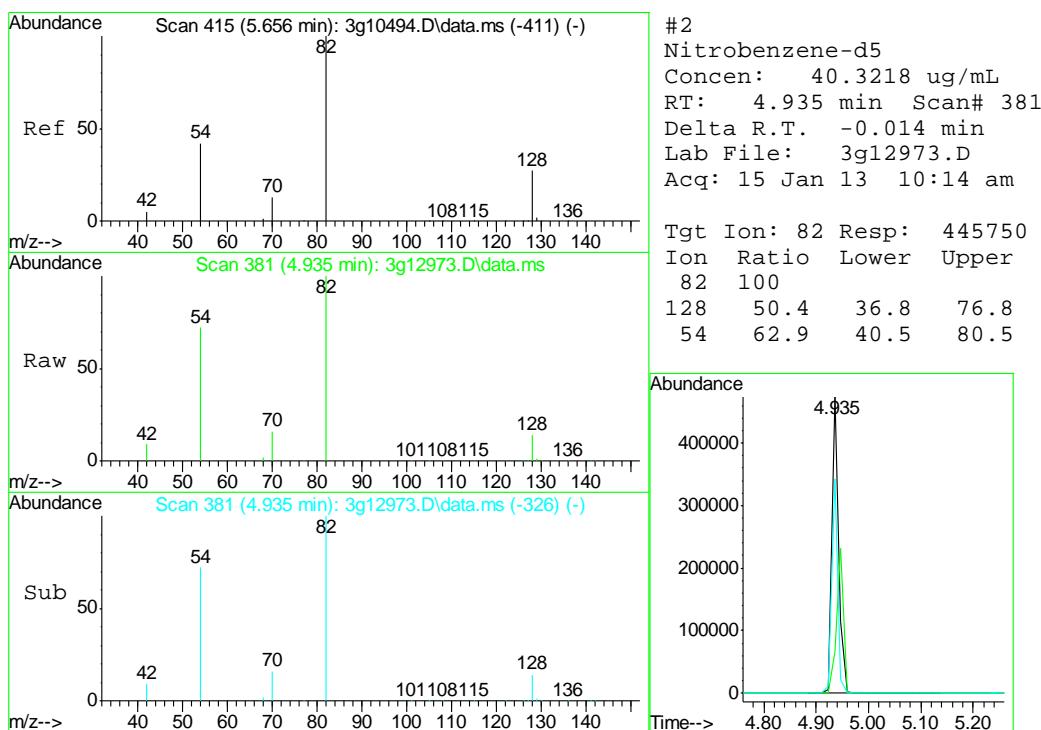
Quant Time: Jan 15 13:46:44 2013  
 Quant Method : C:\msdchem\1\METHODS\SIMPE3G611.M  
 Quant Title : PAHSIM BASE  
 QLast Update : Thu Jan 10 14:18:35 2013  
 Response via : Initial Calibration

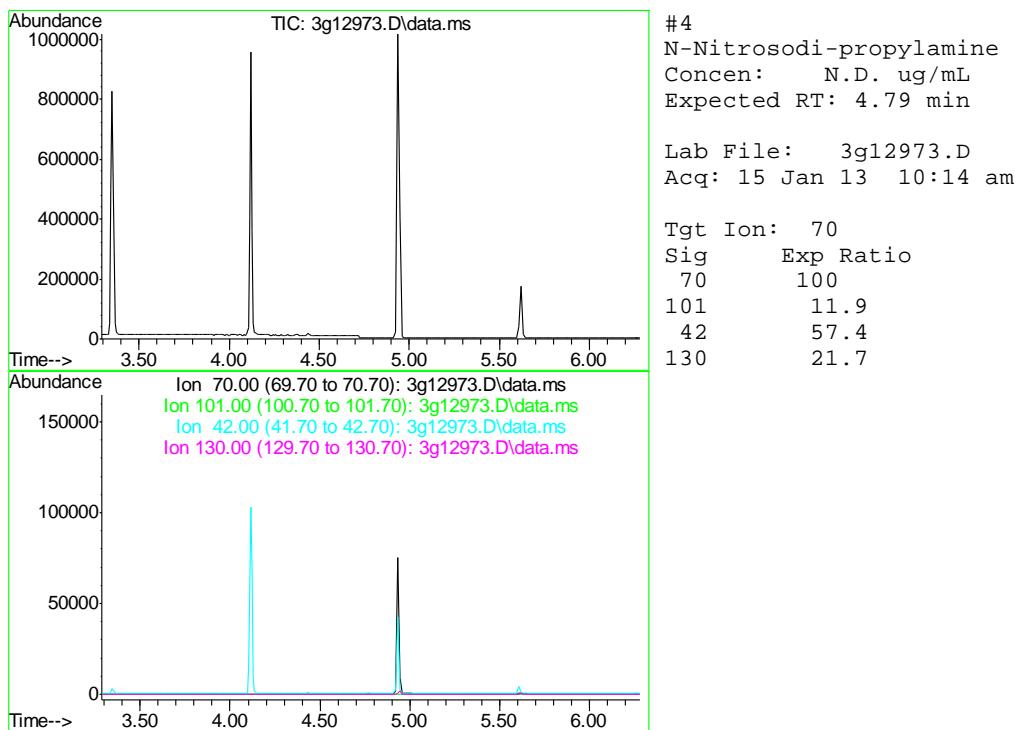
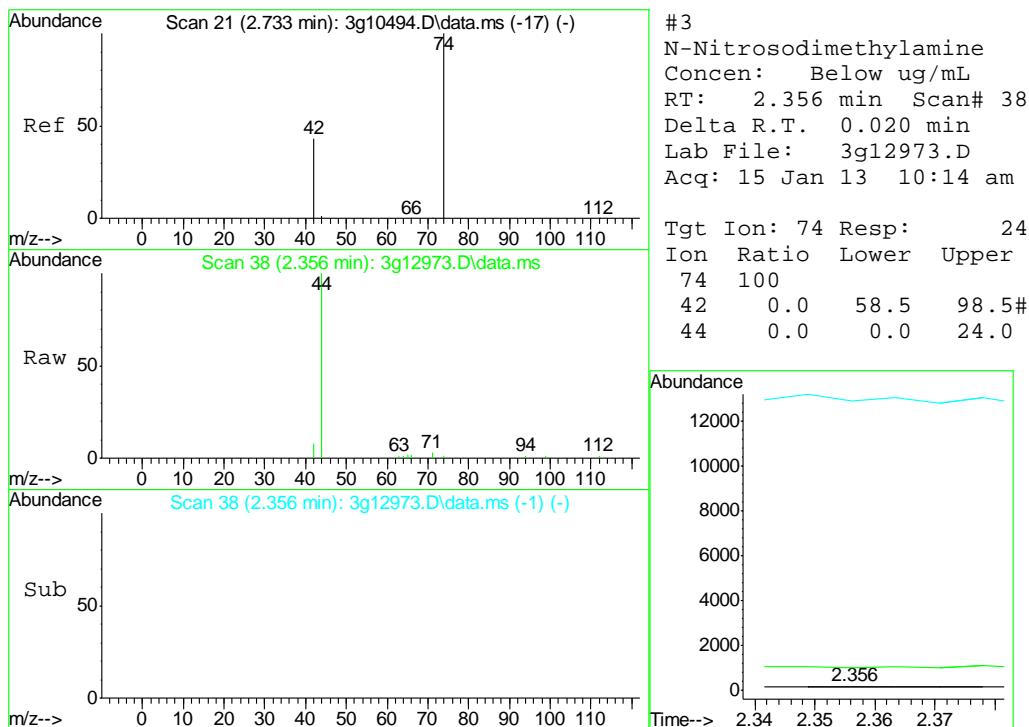


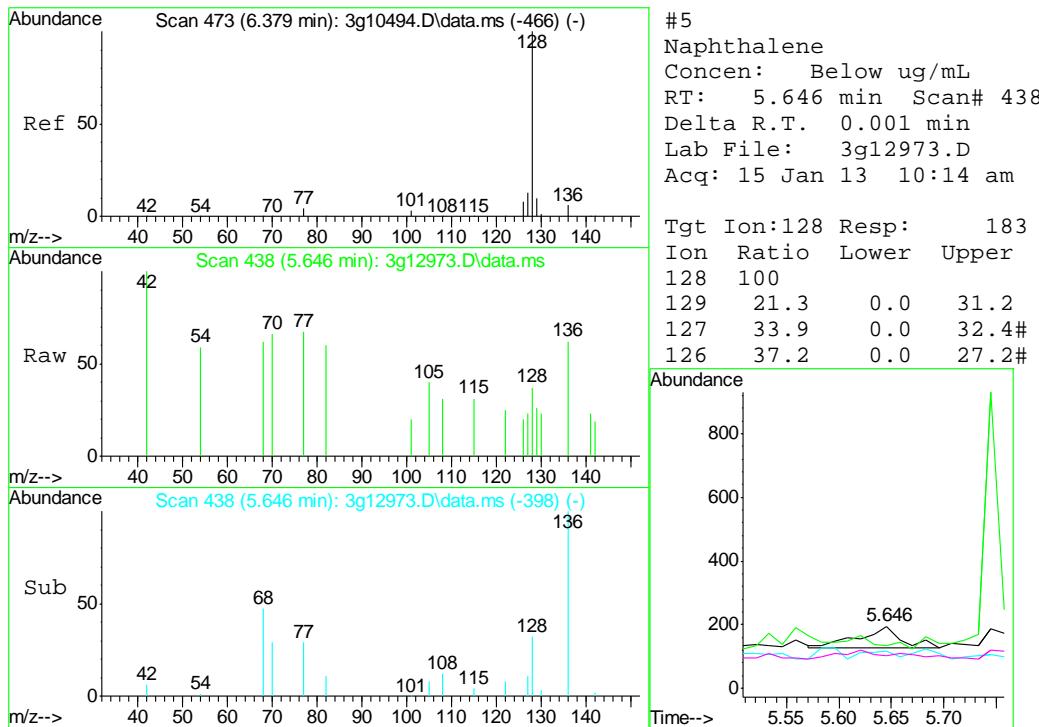


9.2.1

9

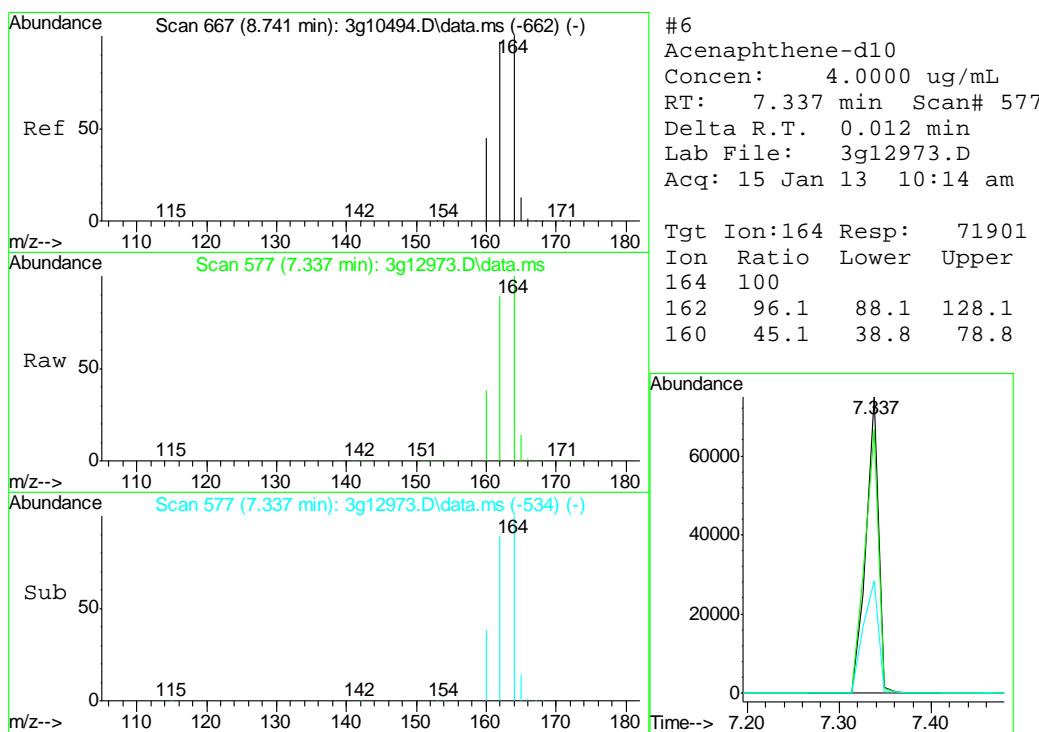


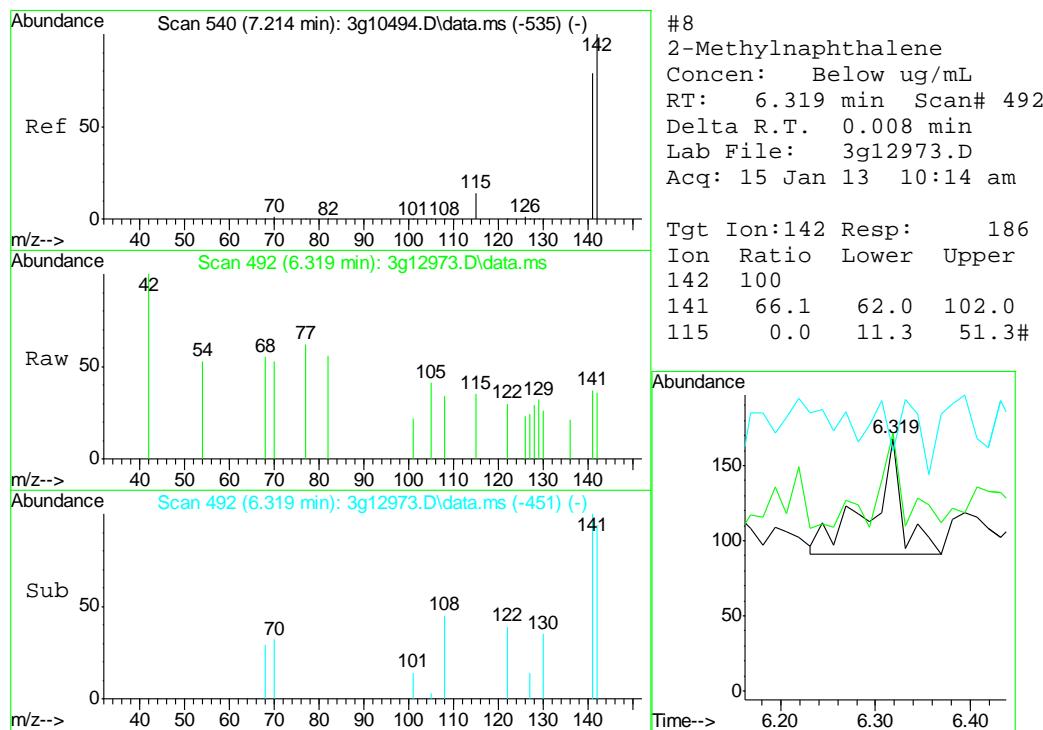
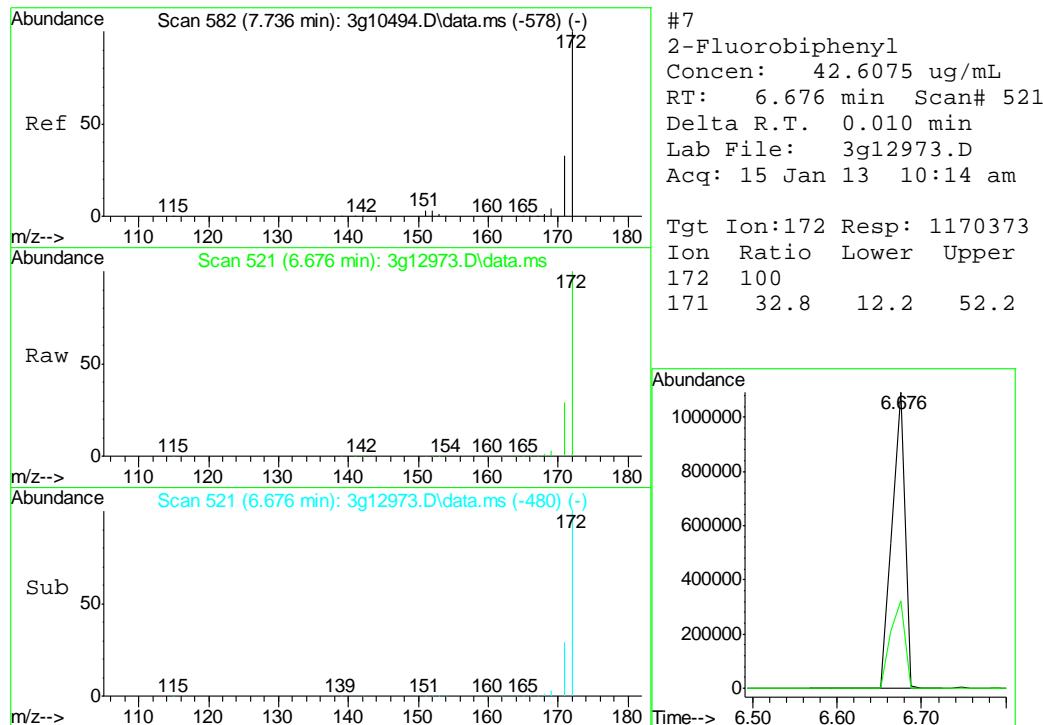


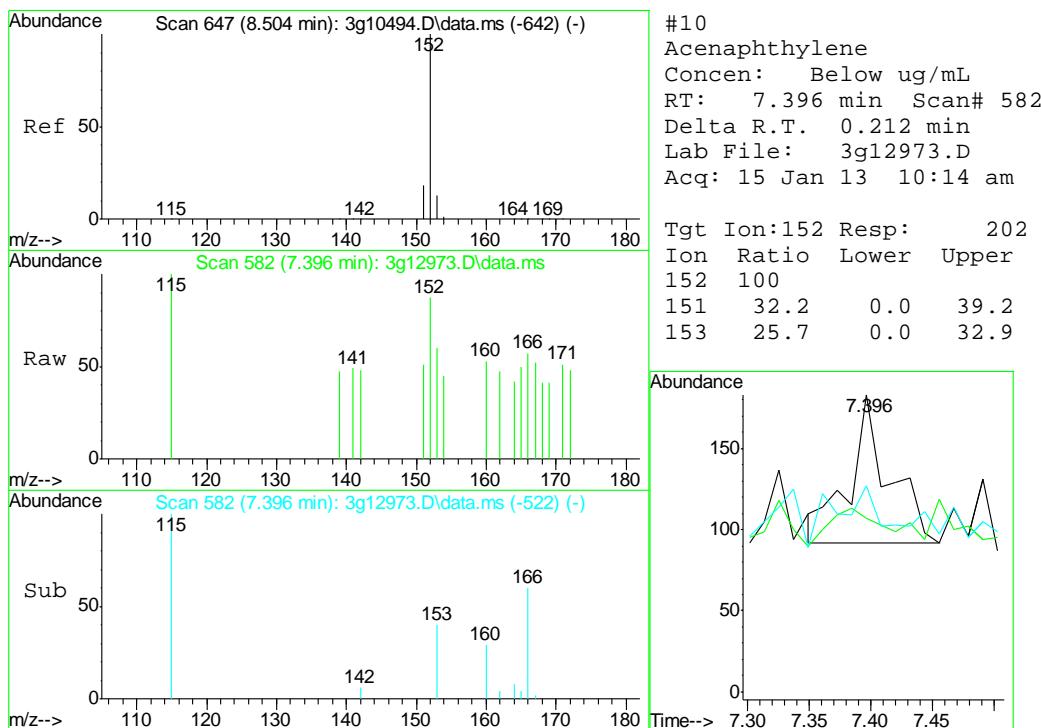
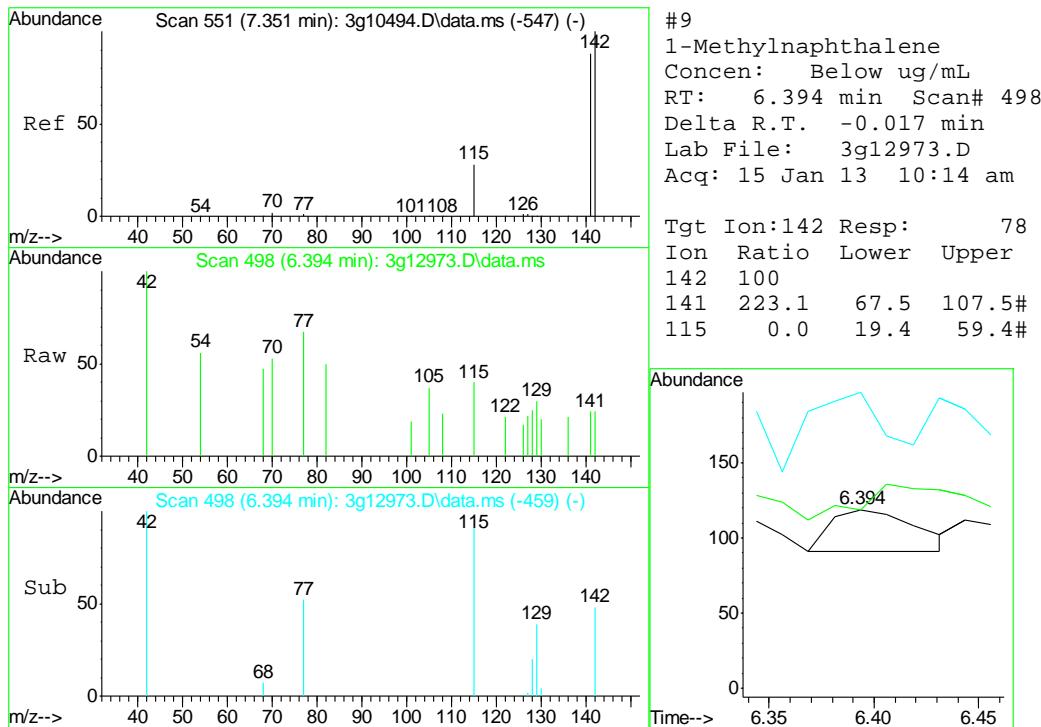


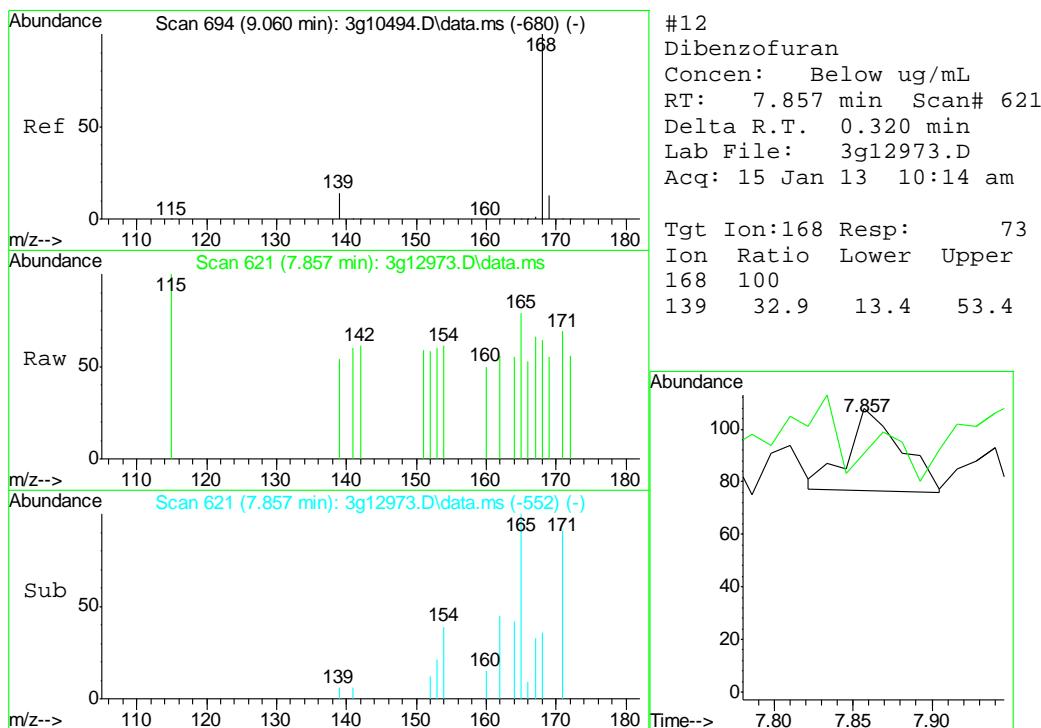
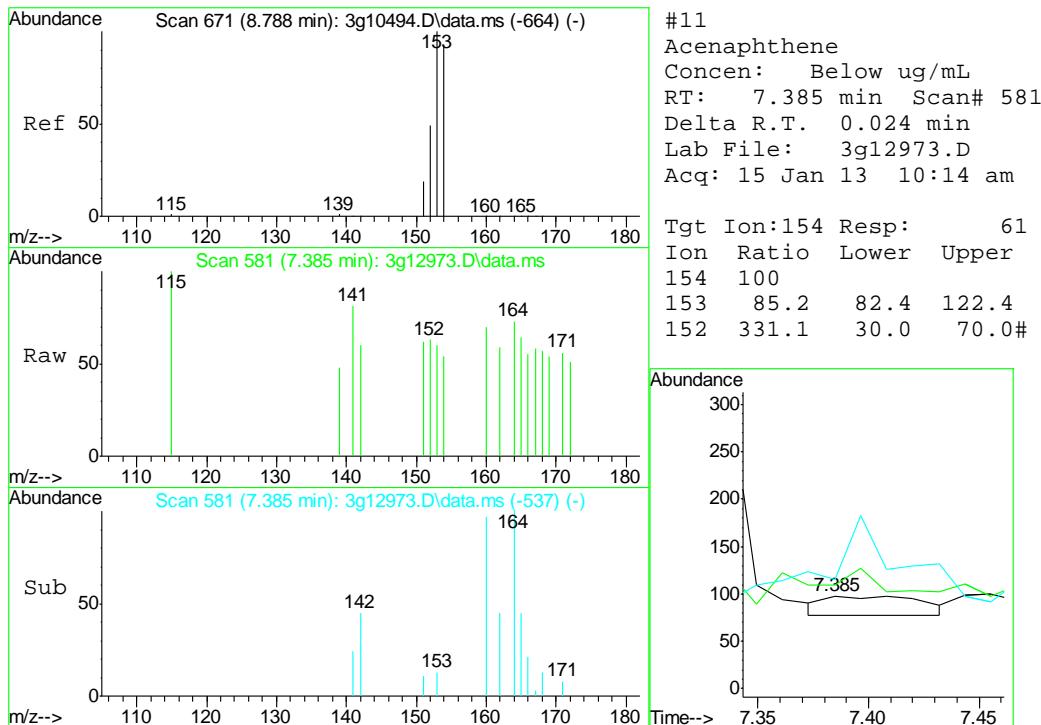
9.2.1

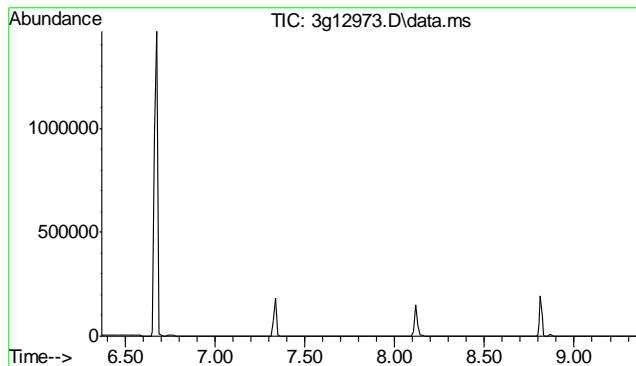
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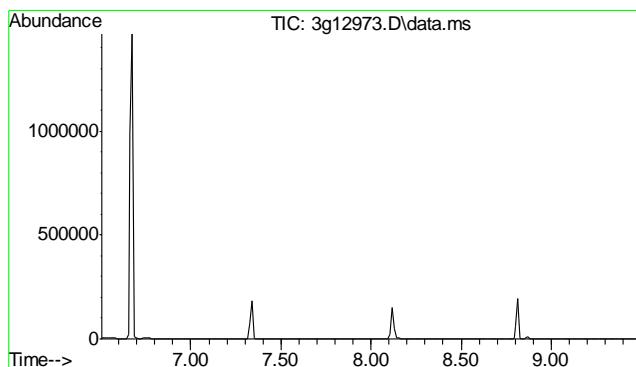
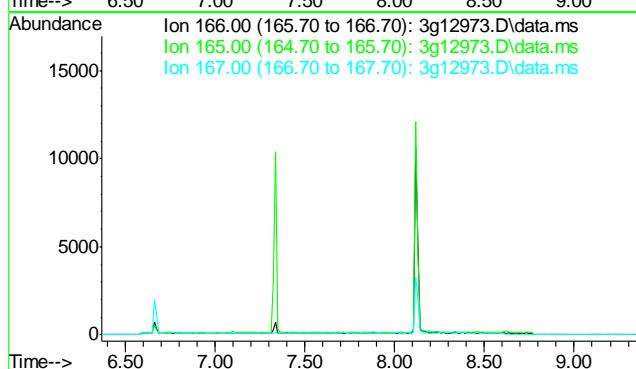




#13  
Fluorene  
Concen: N.D. ug/mL  
Expected RT: 7.87 min

Lab File: 3g12973.D  
Acq: 15 Jan 13 10:14 am

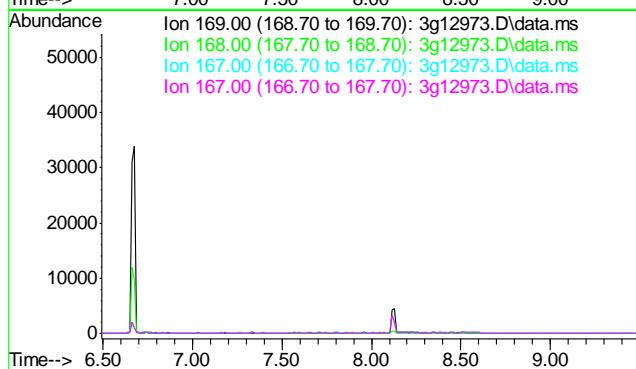
Tgt Ion:	166
Sig	Exp Ratio
166	100
165	92.0
167	13.1

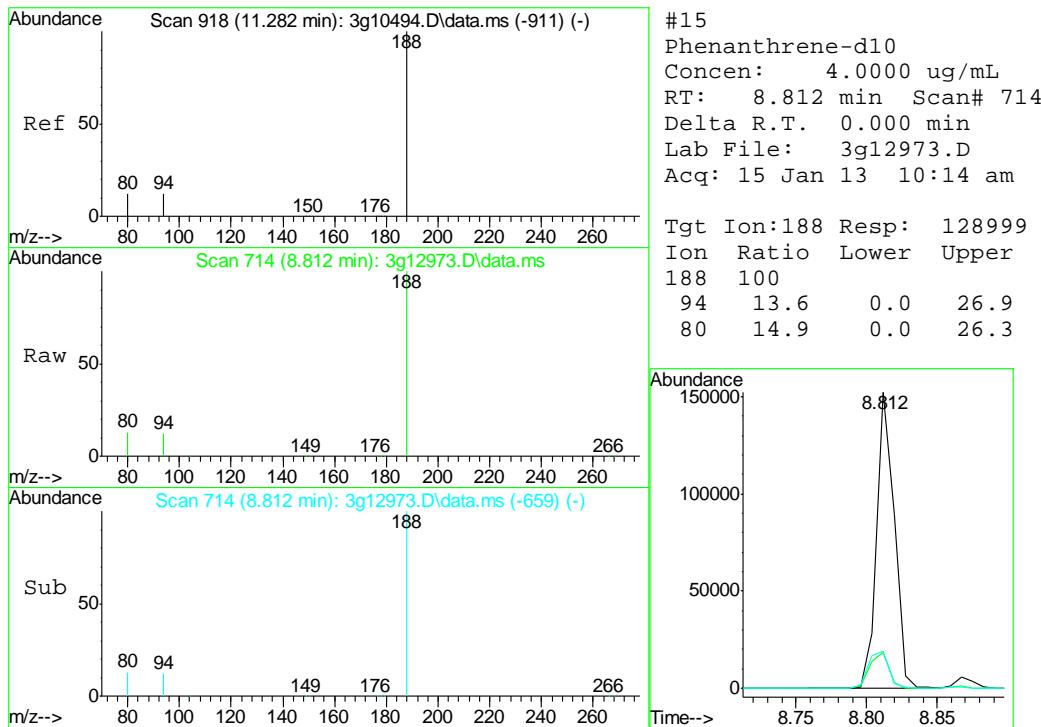


#14  
Diphenylamine  
Concen: N.D. ug/mL  
Expected RT: 8.00 min

Lab File: 3g12973.D  
Acq: 15 Jan 13 10:14 am

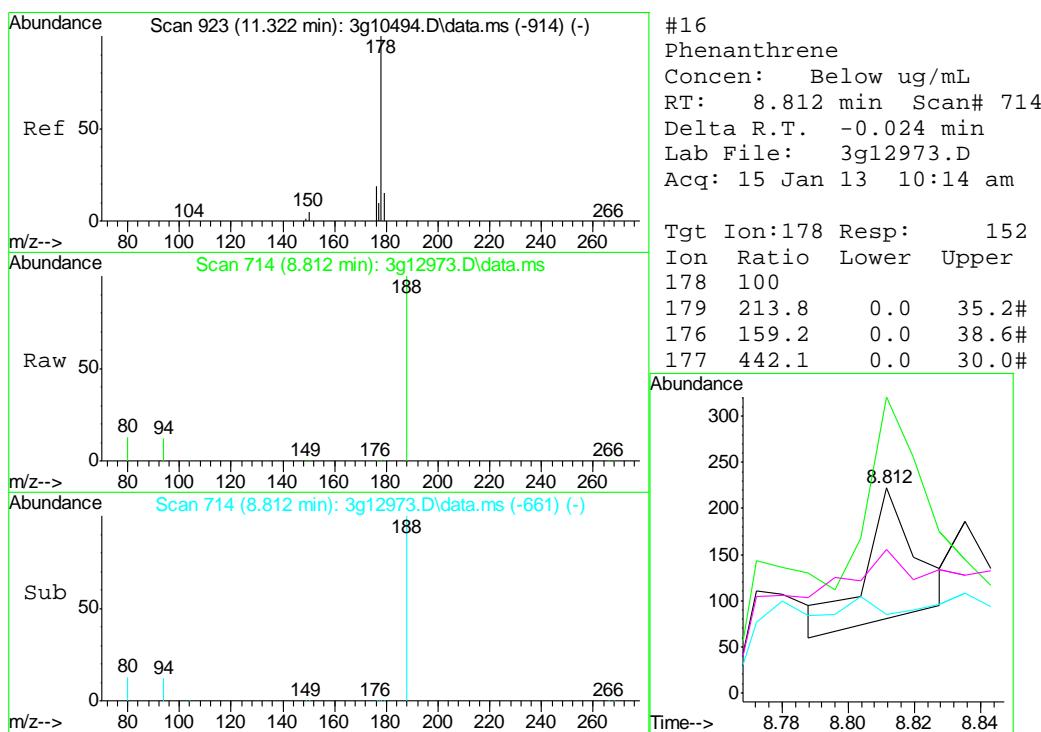
Tgt Ion:	169
Sig	Exp Ratio
169	100
168	61.7
167	34.1
167	34.1

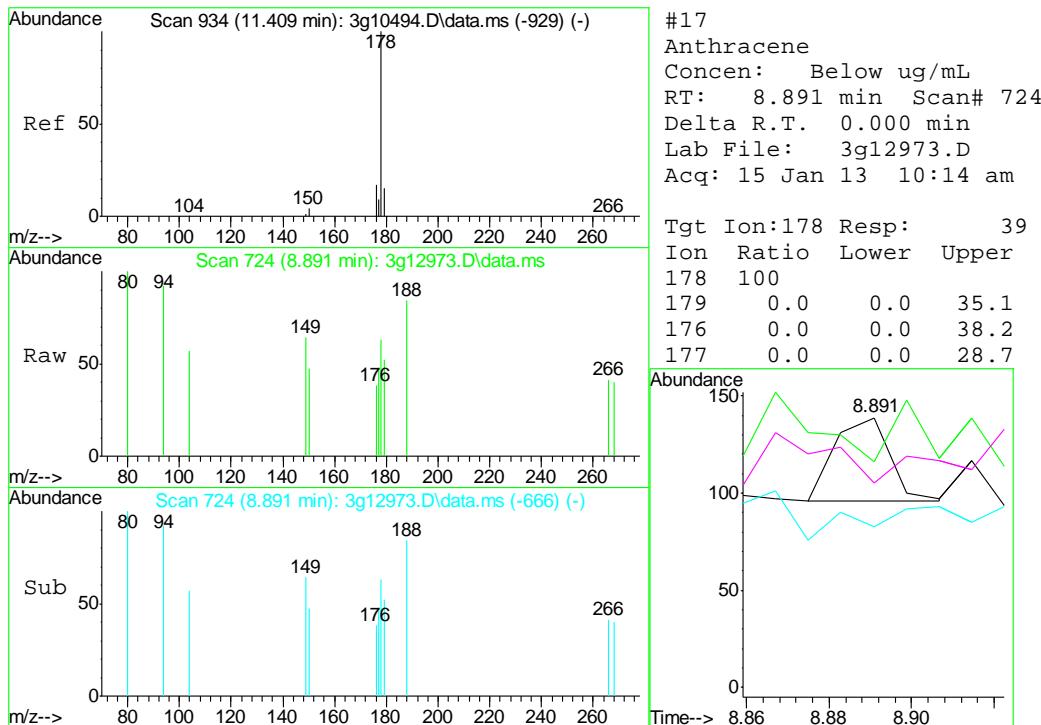




9.2.1

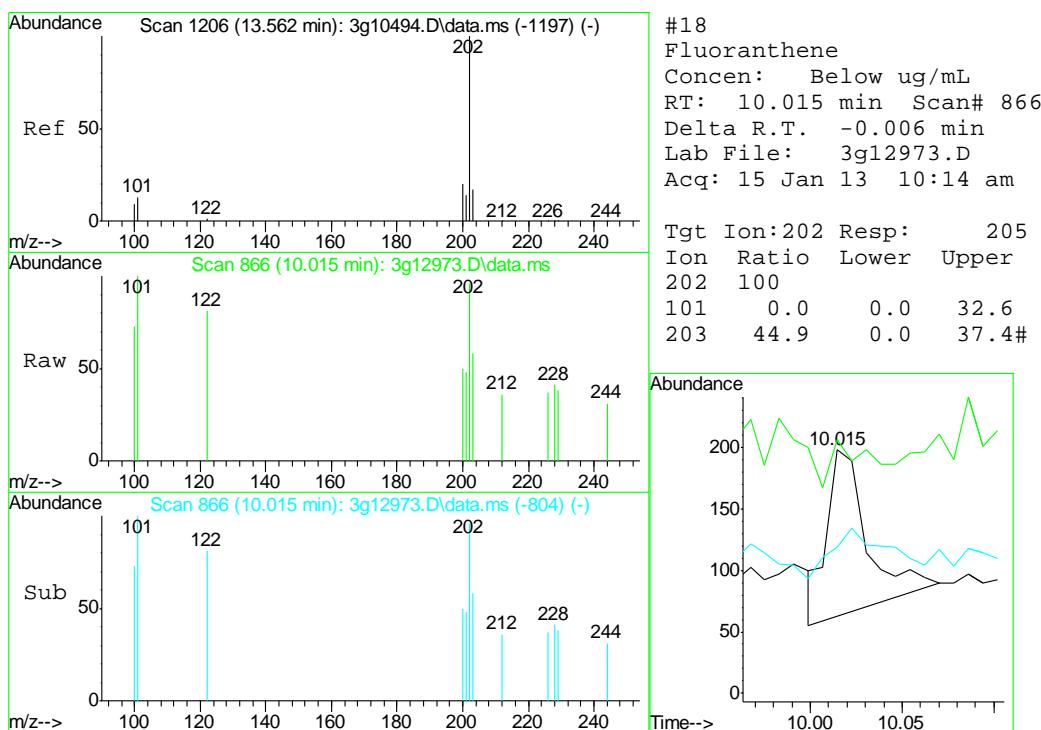
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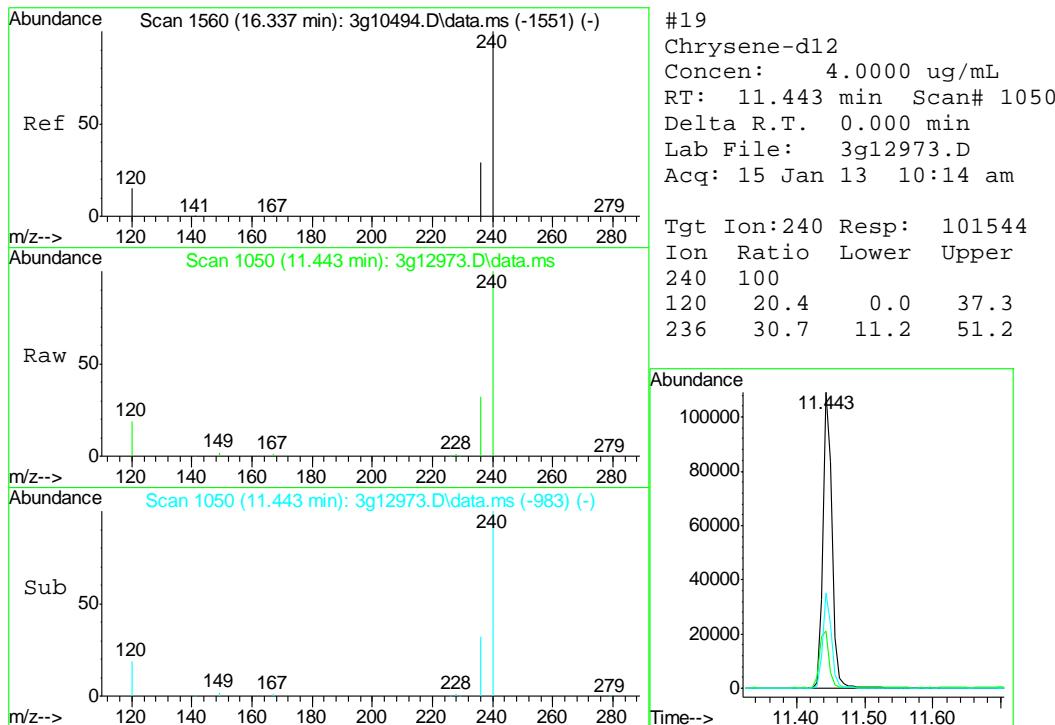




9.2.1

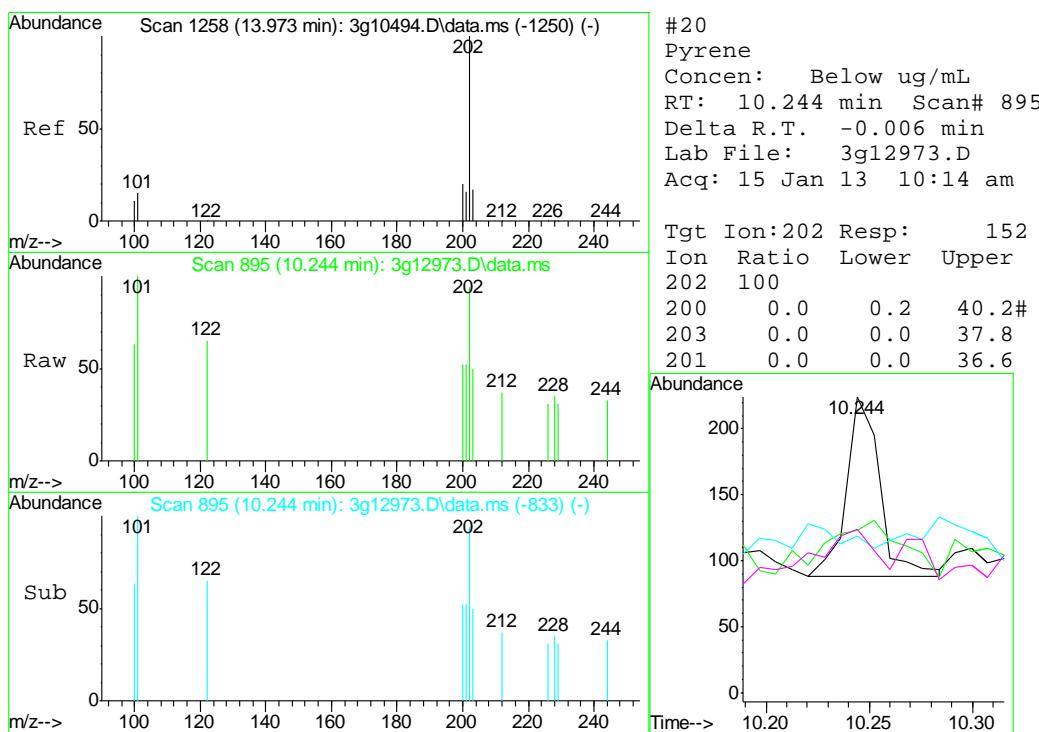
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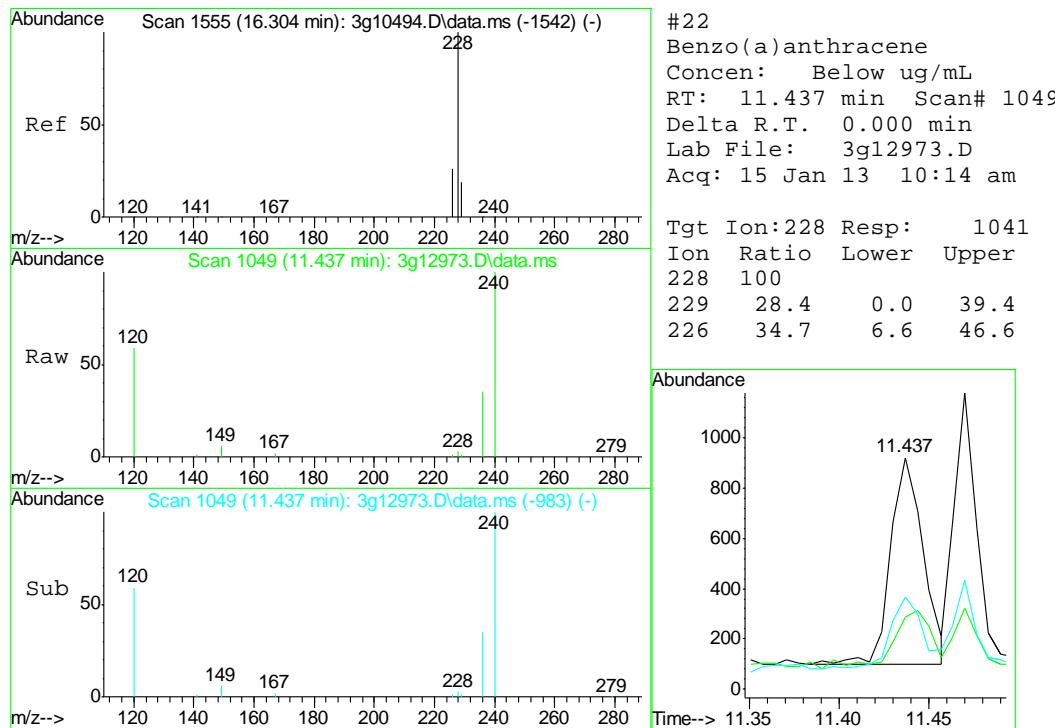
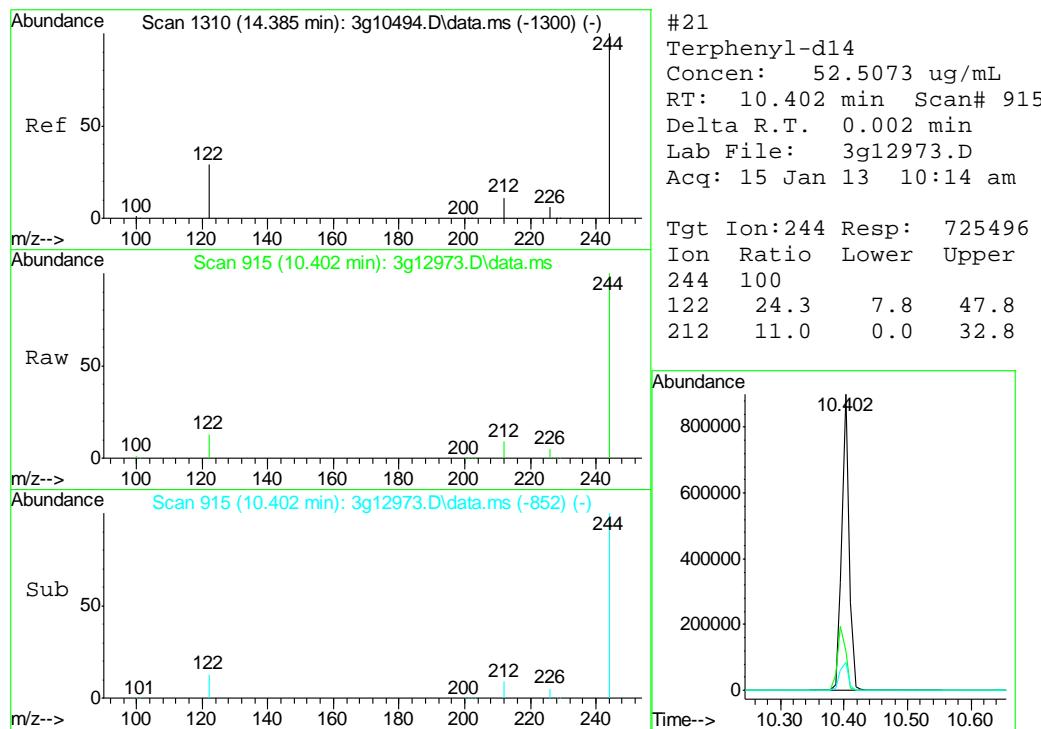


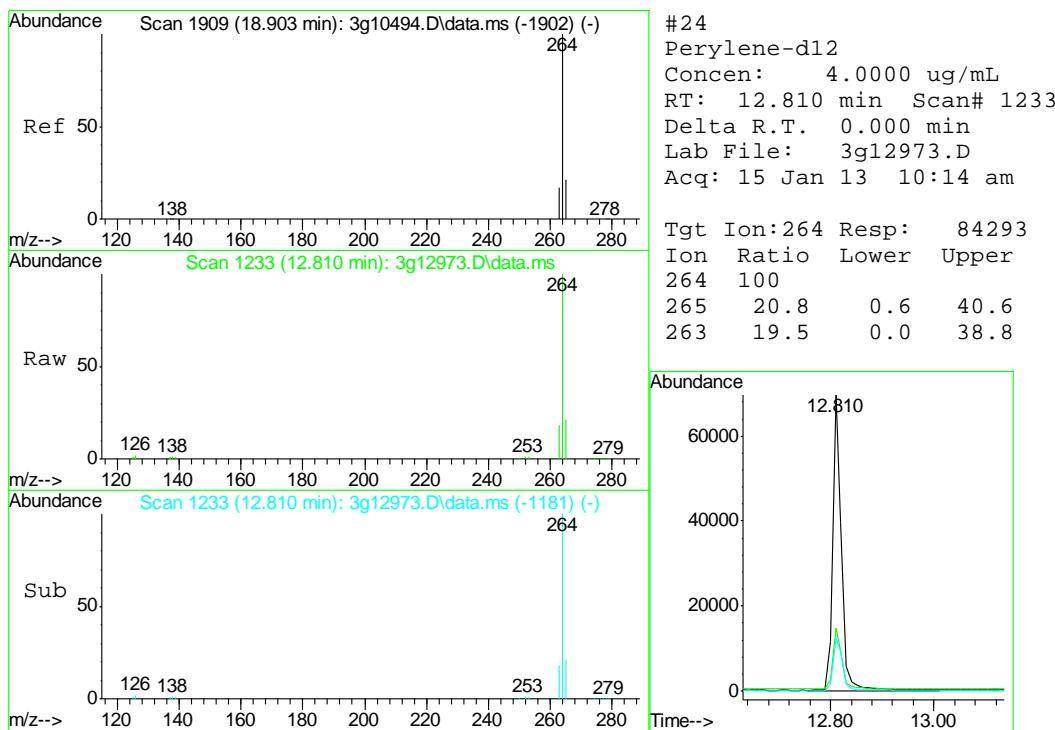
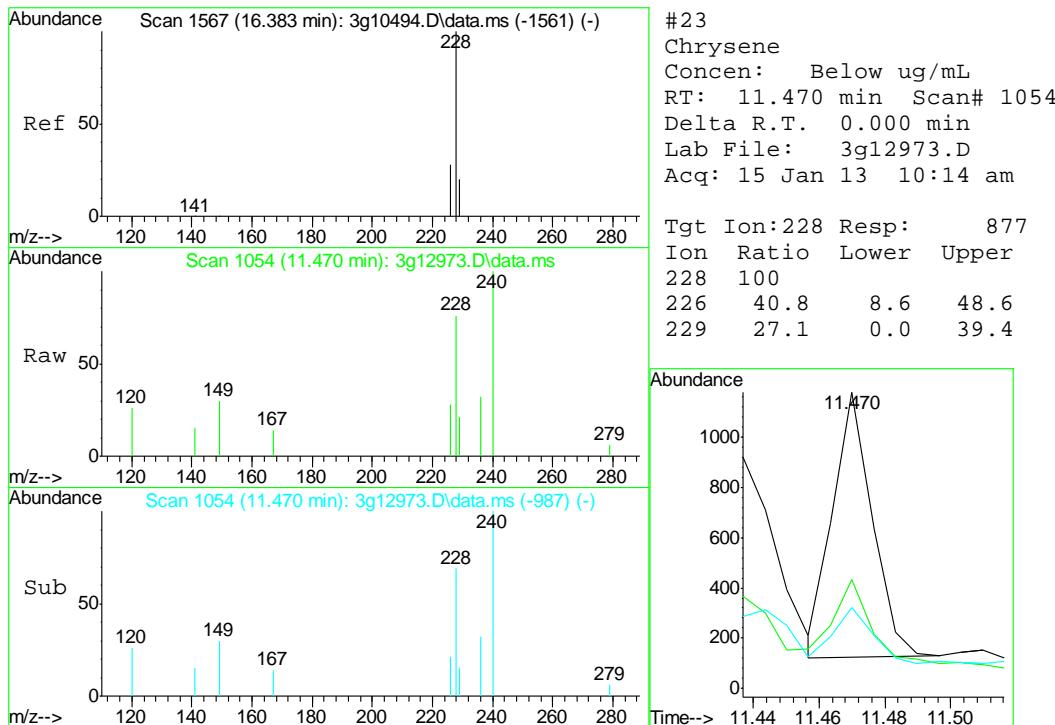


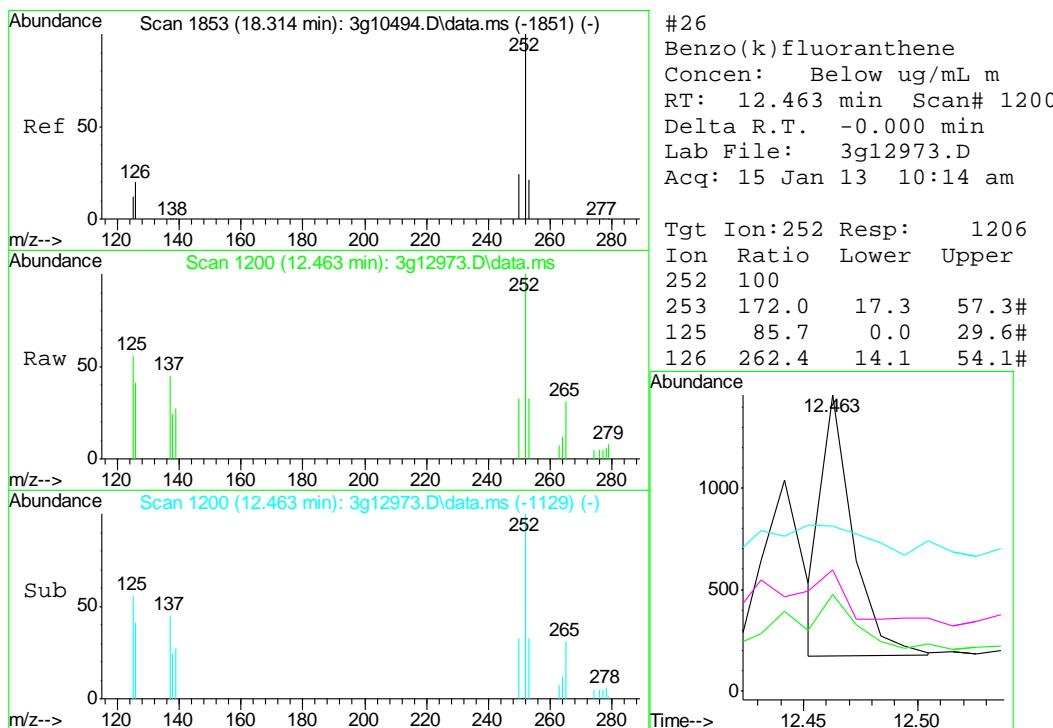
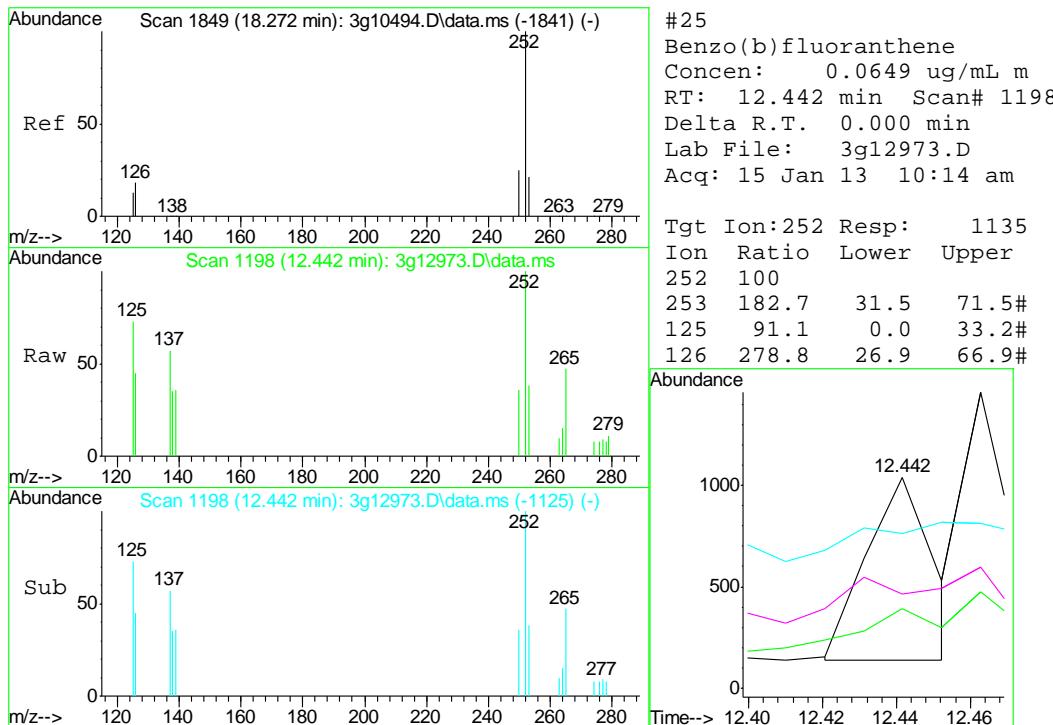
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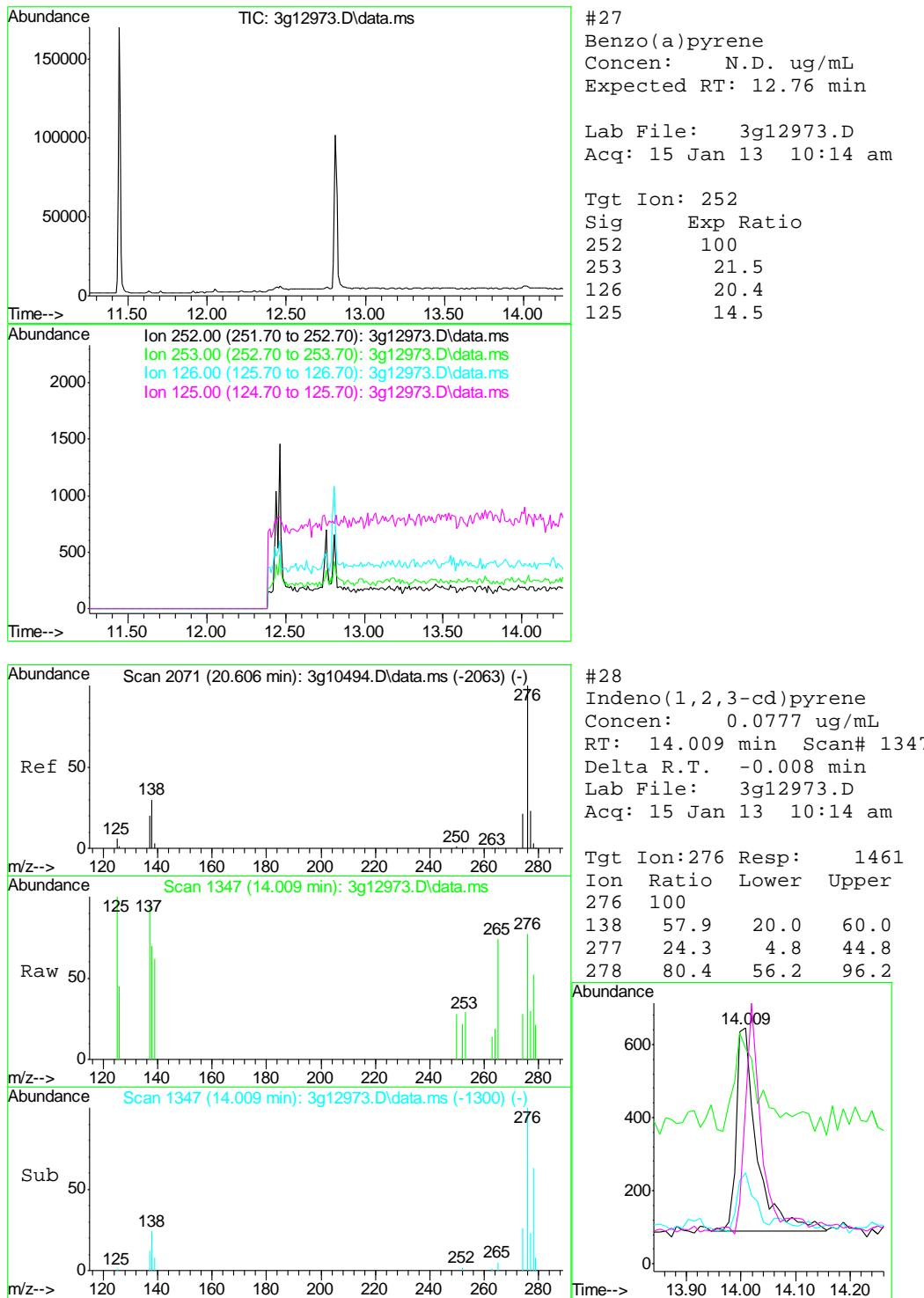
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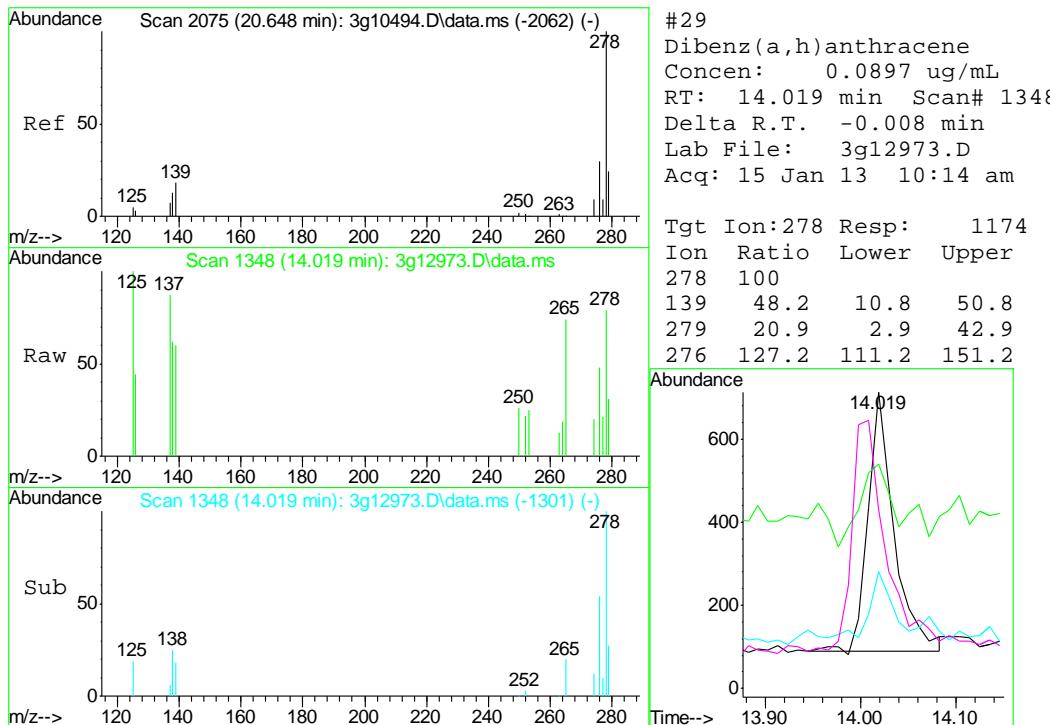






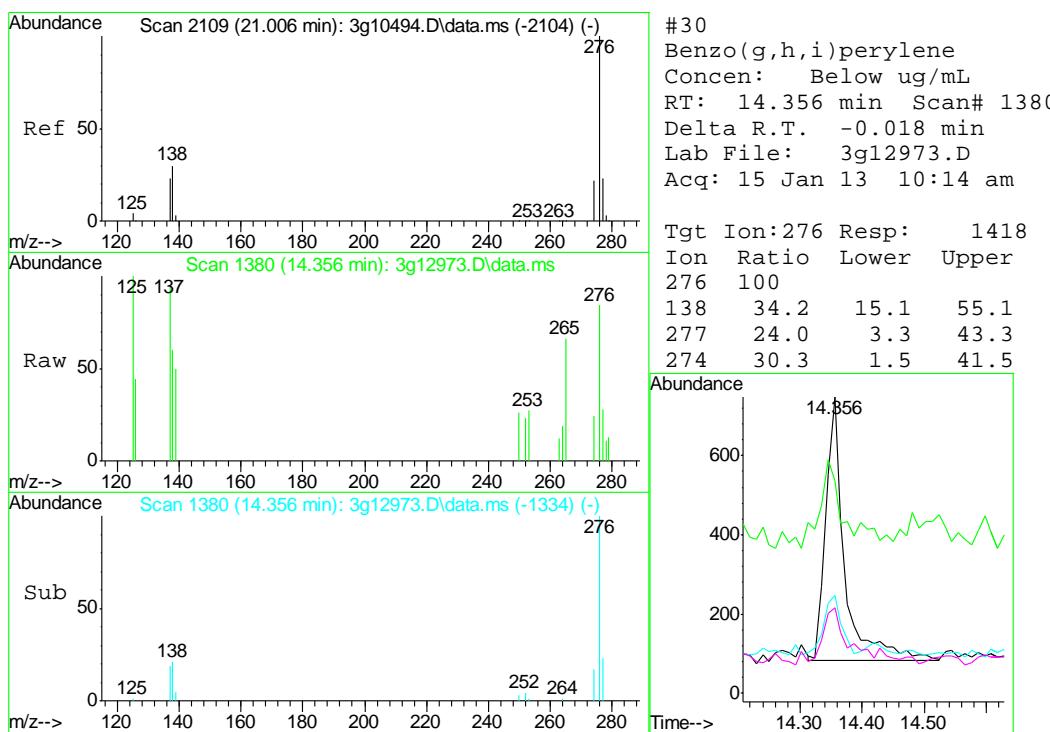






9.2.1

9





## GC Volatiles

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### QC Data Summaries

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Includes the following where applicable:

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

**Method Blank Summary**

**Job Number:** D42556  
**Account:** XTOKWR XTO Energy  
**Project:** PCU 296-5A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
GGB1045-MB	GB19116.D	1	01/11/13	SK	n/a	n/a	GGB1045

The QC reported here applies to the following samples:

**Method:** SW846 8015B

D42556-1, D42556-2

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	90% 60-140%

10.1.1

10

## Blank Spike Summary

Page 1 of 1

Job Number: D42556

Account: XTOKWR XTO Energy

Project: PCU 296-5A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
GGB1045-BS	GB19117.D	1	01/11/13	SK	n/a	n/a	GGB1045

The QC reported here applies to the following samples:

Method: SW846 8015B

D42556-1, D42556-2

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**

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\* = Outside of Control Limits.

# Matrix Spike/Matrix Spike Duplicate Summary

Page 1 of 1

Job Number: D42556

Account: XTOKWR XTO Energy

Project: PCU 296-5A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
D42466-11MS	GB19119.D	1	01/11/13	SK	n/a	n/a	GGB1045
D42466-11MSD	GB19120.D	1	01/11/13	SK	n/a	n/a	GGB1045
D42466-11	GB19118.D	1	01/11/13	SK	n/a	n/a	GGB1045

The QC reported here applies to the following samples:

Method: SW846 8015B

D42556-1, D42556-2

CAS No.	Compound	D42466-11		Spike	MS	MS	MSD	MSD	RPD	Limits Rec/RPD
		mg/kg	Q	mg/kg	mg/kg	%	mg/kg	%		
	TPH-GRO (C6-C10)	ND		152	170	112	167	110	2	70-130/30

CAS No.	Surrogate Recoveries	MS	MSD	D42466-11	Limits
120-82-1	1,2,4-Trichlorobenzene	109%	100%	84%	60-140%

\* = Outside of Control Limits.

10.3.1  
10



## GC Volatiles

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Raw Data

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**Manual Integrations  
APPROVED  
(compounds with "m" flag)**  
**Judy Nelson  
01/14/13 12:14**

## Quantitation Report (QT Reviewed)

Signal #1 : Y:\1\DATA\011113\GB19132.D\FID1A.CH Vial: 19  
 Signal #2 : Y:\1\DATA\011113\GB19132.D\FID2B.CH  
 Acq On : 11 Jan 2013 9:59 pm Operator: StephK  
 Sample : D42556-1, 50X Inst : GC/MS Ins  
 Misc : GC3352,GGB1045,5.066,,100,5,1 Multiplr: 1.00  
 IntFile Signal #1: TVH1.E IntFile Signal #2: FB2.E  
 Quant Time: Jan 14 08:43:57 2013 Quant Results File: TB868GB868SOIL.RES

Quant Method : C:\MSDCHEM\1...\TB868GB868SOIL.M (Chemstation Integrator)  
 Title : 8015B/8021B TVH/BTEX  
 Last Update : Fri Jan 11 14:33:28 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
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**System Monitoring Compounds**

2) S	1,2,4-Trichlorobenzene	14.36	2905471	92.726 %	m
10) S	1,2,4-Trichlorobenzene (P)	14.36	14820177	91.186 %	m

**Target Compounds**

1) H	TVH-Gasoline	7.23	3738403	<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	7.66	104796	0.264	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	0.00	0	N.D.	ug/L d

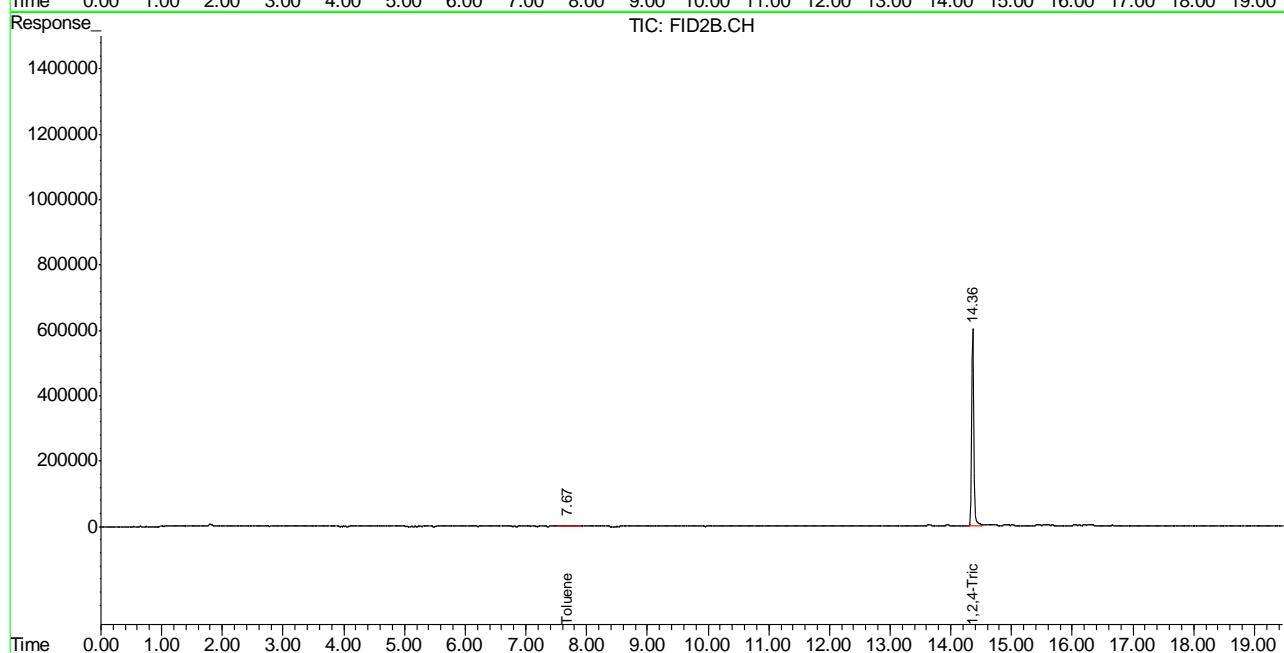
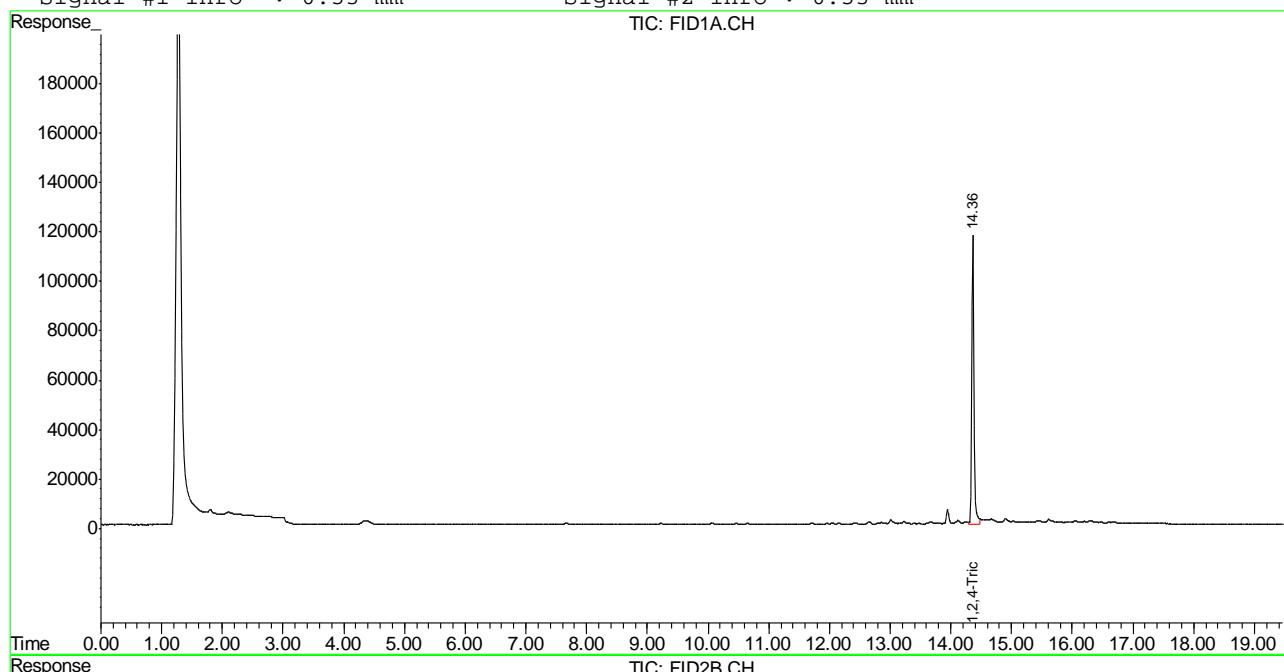
-----  
 (f)=RT Delta > 1/2 Window (m)=manual int.  
 GB19132.D TB868GB868SOIL.M Mon Jan 14 09:01:25 2013 GC

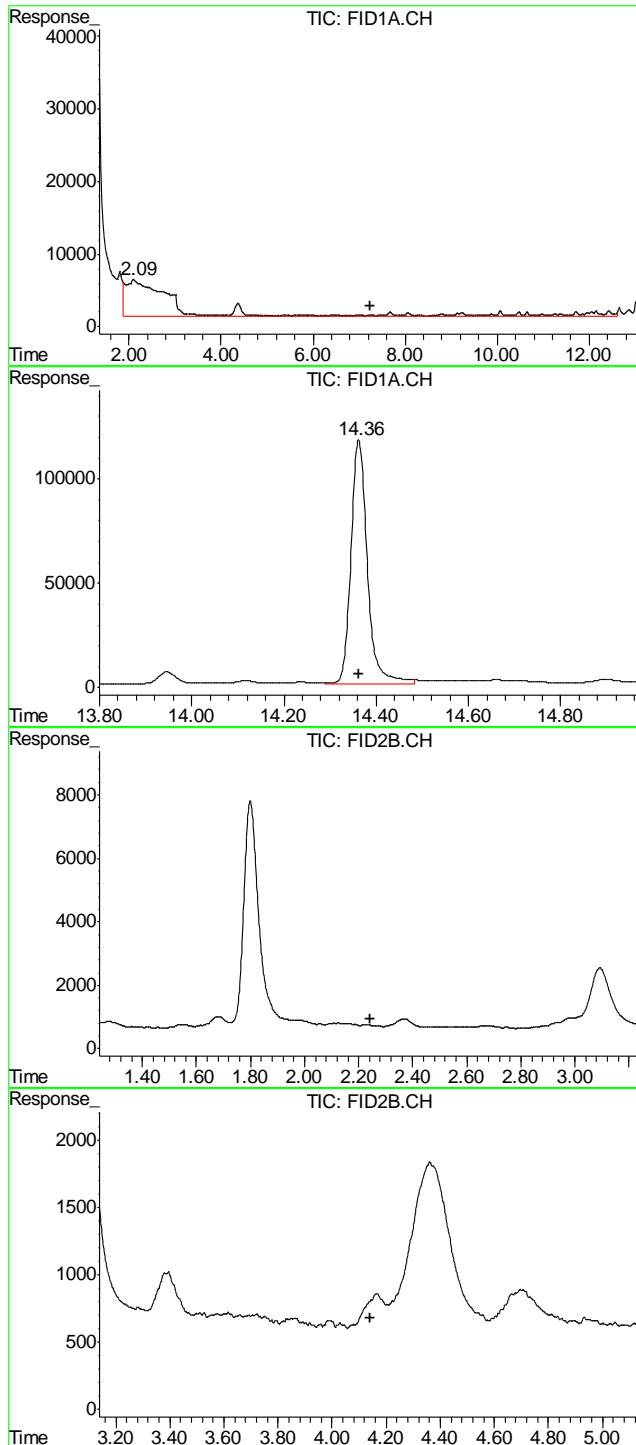
## Quantitation Report (QT Reviewed)

Signal #1 : Y:\1\DATA\011113\GB19132.D\FID1A.CH Vial: 19  
 Signal #2 : Y:\1\DATA\011113\GB19132.D\FID2B.CH  
 Acq On : 11 Jan 2013 9:59 pm Operator: StephK  
 Sample : D42556-1, 50X Inst : GC/MS Ins  
 Misc : GC3352,GGB1045,5.066,,100,5,1 Multiplr: 1.00  
 IntFile Signal #1: TVH1.E IntFile Signal #2: FB2.E  
 Quant Time: Jan 14 8:56 2013 Quant Results File: TB868GB868SOIL.RES

Quant Method : C:\MSDCHEM\1...\TB868GB868SOIL.M (Chemstation Integrator)  
 Title : 8015B/8021B TVH/BTEX  
 Last Update : Fri Jan 11 14:33:28 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: 3738403  
 Conc: N.D.

## #2 1,2,4-Trichlorobenzene

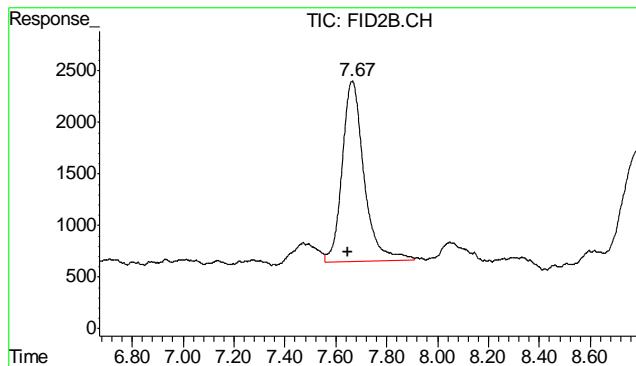
R.T.: 14.362 min  
 Delta R.T.: 0.000 min  
 Response: 2905471  
 Conc: 92.73 % m

## #4 Methyl-t-butyl-ether

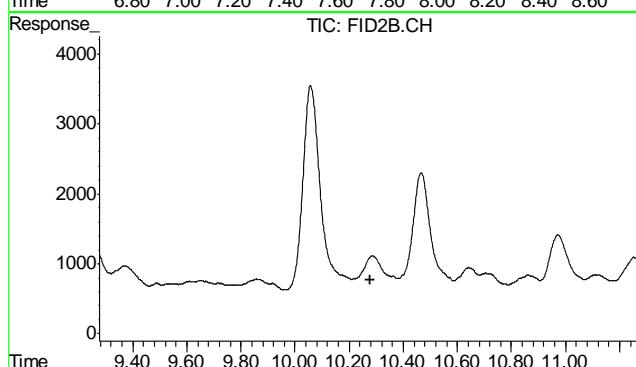
R.T.: 0.000 min  
 Exp R.T. : 2.241 min  
 Response: 0  
 Conc: N.D.

## #5 Benzene

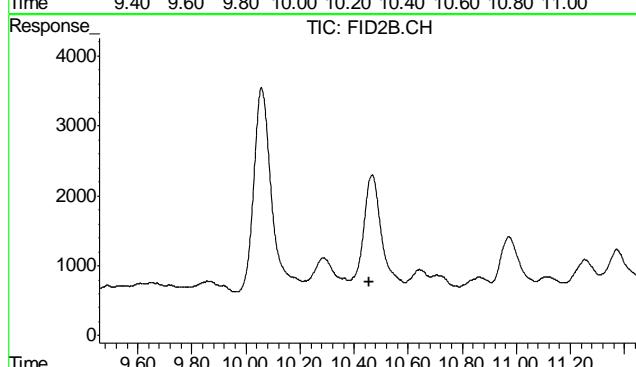
R.T.: 0.000 min  
 Exp R.T. : 4.139 min  
 Response: 0  
 Conc: N.D.



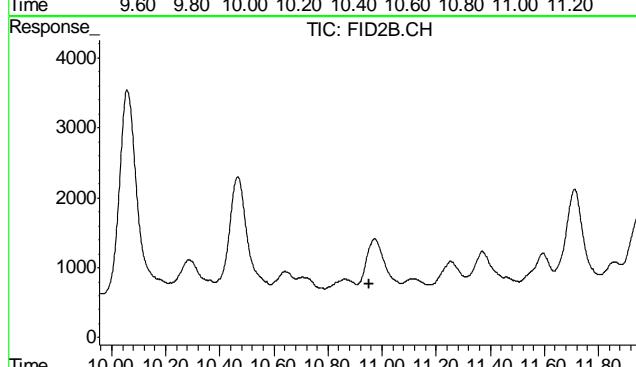
#6 Toluene  
R.T.: 7.665 min  
Delta R.T.: 0.017 min  
Response: 104796  
Conc: 0.26 ug/L



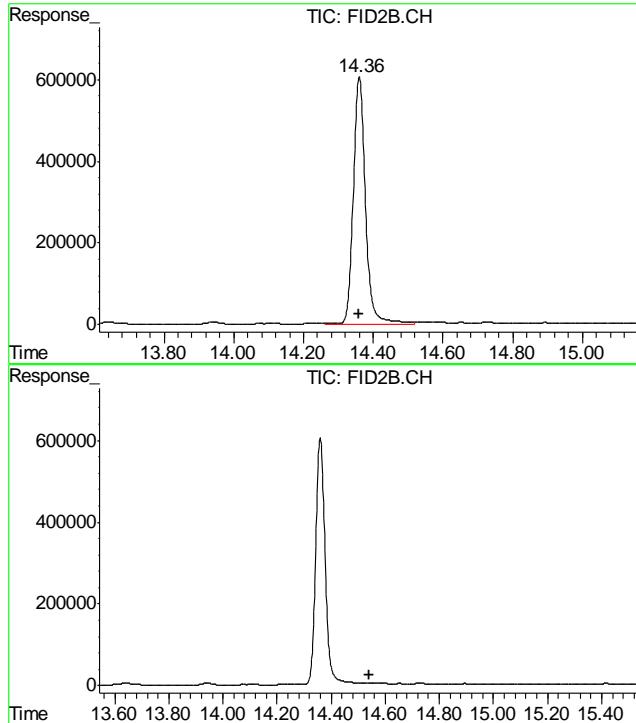
#7 Ethylbenzene  
R.T.: 0.000 min  
Exp R.T. : 10.276 min  
Response: 0  
Conc: N.D.



#8 m,p-Xylene  
R.T.: 0.000 min  
Exp R.T. : 10.457 min  
Response: 0  
Conc: N.D.



#9 o-Xylene  
R.T.: 0.000 min  
Exp R.T. : 10.954 min  
Response: 0  
Conc: N.D.



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

R.T.: 14.359 min

Delta R.T.: 0.000 min

Response: 14820177

Conc: 91.19 % m

#11 Naphthalene

R.T.: 0.000 min

Exp R.T. : 14.542 min

Response: 0

Conc: N.D.

11.1.1

Manual Integrations  
APPROVED  
(compounds with "m" flag)

Judy Nelson  
01/14/13 12:14

## Quantitation Report (QT Reviewed)

Signal #1 : Y:\1\DATA\011113\GB19133.D\FID1A.CH Vial: 20  
 Signal #2 : Y:\1\DATA\011113\GB19133.D\FID2B.CH  
 Acq On : 11 Jan 2013 10:35 pm Operator: StephK  
 Sample : D42556-2, 50X Inst : GC/MS Ins  
 Misc : GC3352,GGB1045,5.060,,100,5,1 Multiplr: 1.00  
 IntFile Signal #1: TVH1.E IntFile Signal #2: FB2.E  
 Quant Time: Jan 14 08:44:01 2013 Quant Results File: TB868GB868SOIL.RES

Quant Method : C:\MSDCHEM\1...\TB868GB868SOIL.M (Chemstation Integrator)  
 Title : 8015B/8021B TVH/BTEX  
 Last Update : Fri Jan 11 14:33:28 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
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## System Monitoring Compounds

2) S	1,2,4-Trichlorobenzene	14.36	2677351	85.446 %	m
10) S	1,2,4-Trichlorobenzene (P)	14.36	13368492	82.254 %	m

## Target Compounds

1) H	TVH-Gasoline	7.23	23992846	0.348	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	7.66	2179890	5.501	ug/L
7) T	Ethylbenzene	10.28	888680	2.627	ug/L m
8) T	m,p-Xylene	10.46	4027842	10.662	ug/L
9) T	o-Xylene	10.96	334429	1.018	ug/L m
11) T	Naphthalene	14.55	226539	1.148	ug/L m

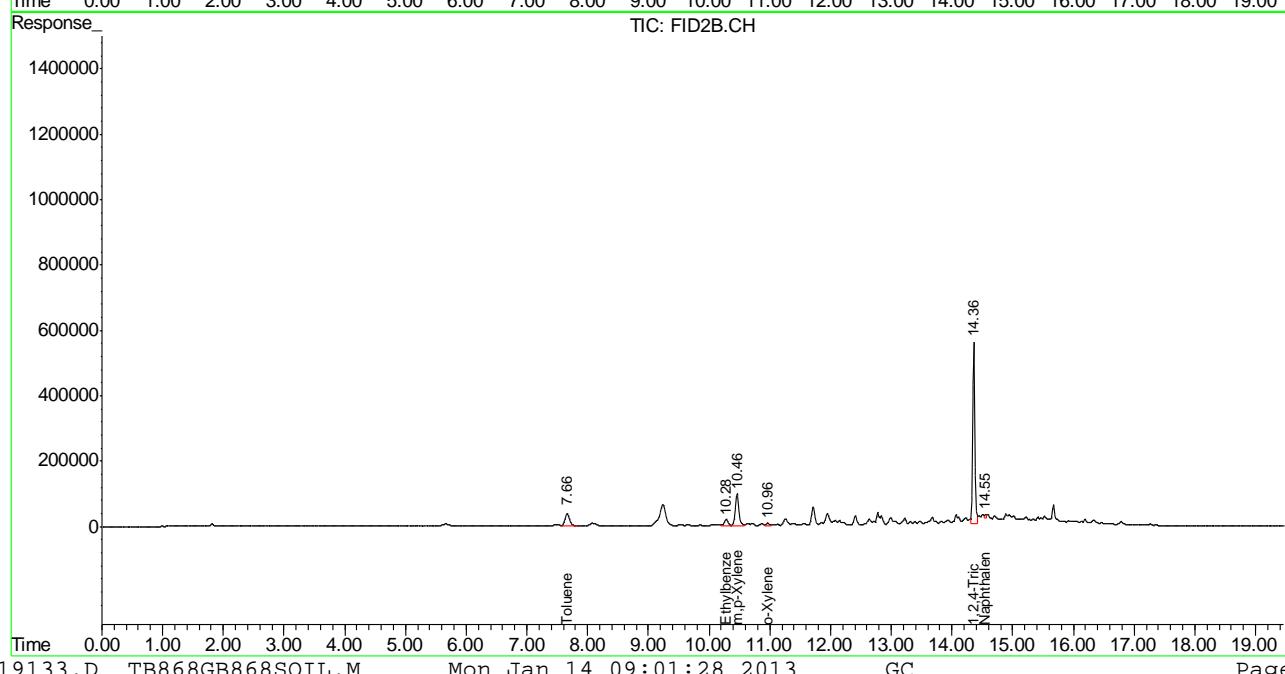
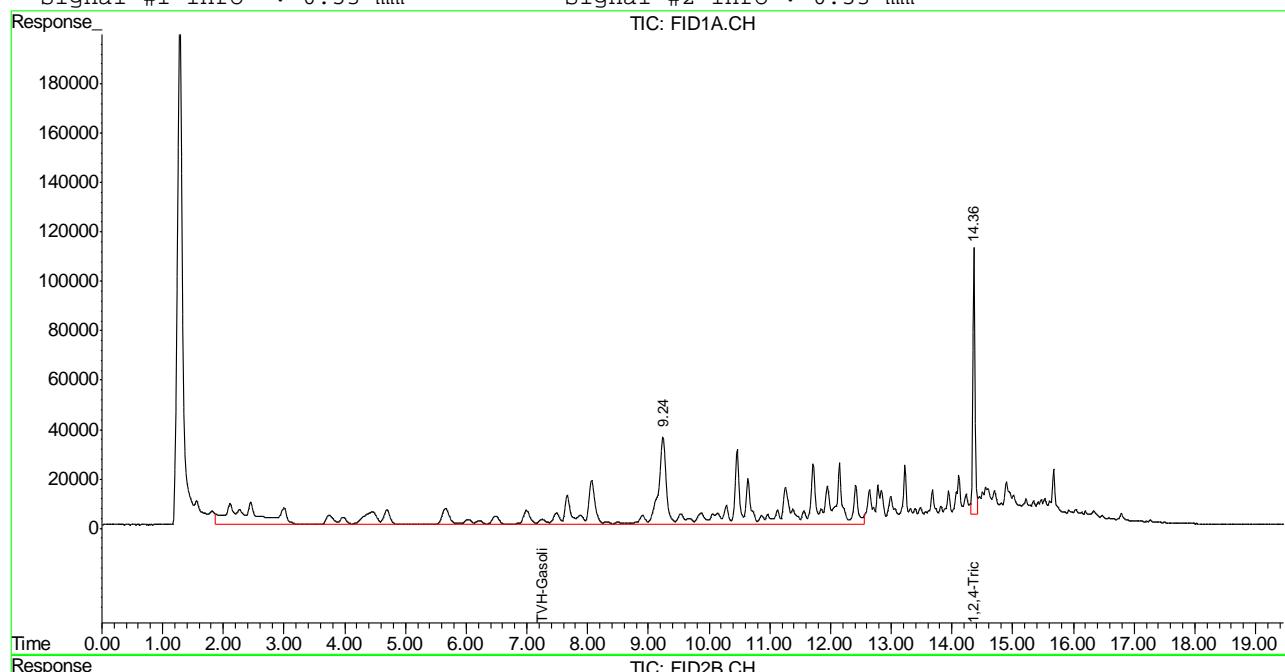
(f)=RT Delta > 1/2 Window (m)=manual int.  
 GB19133.D TB868GB868SOIL.M Mon Jan 14 09:01:28 2013 GC

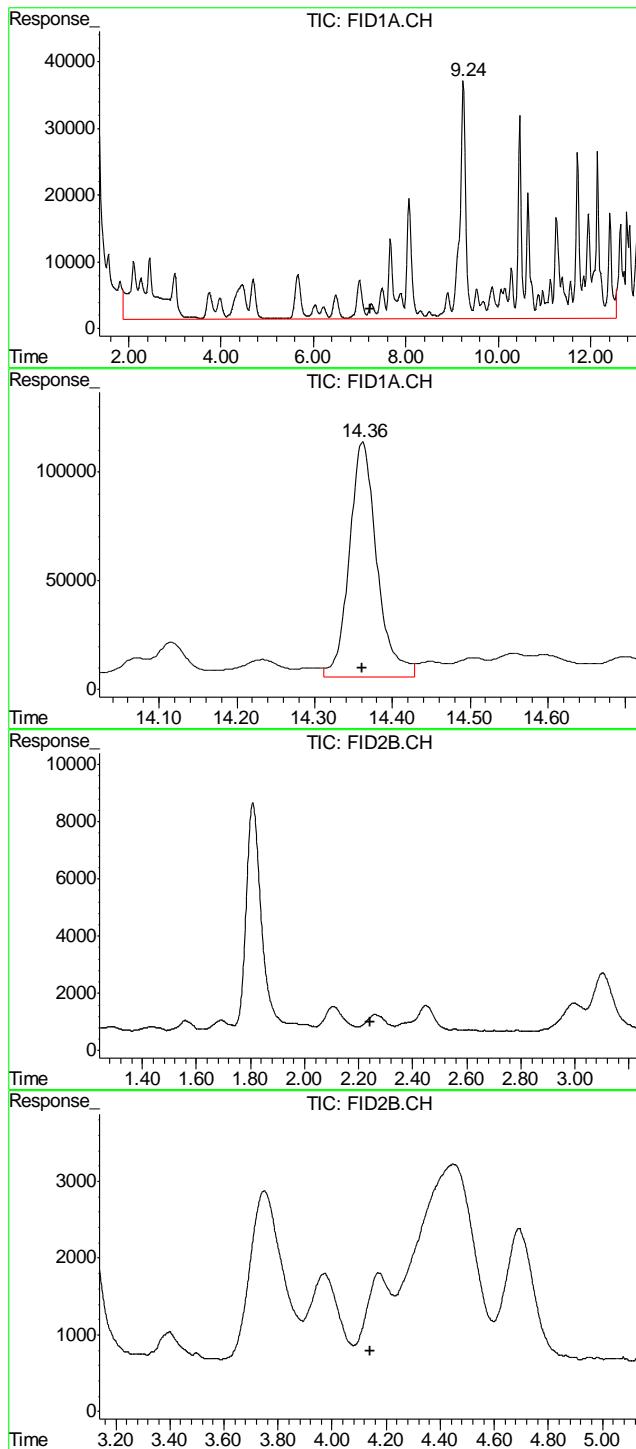
## Quantitation Report (QT Reviewed)

Signal #1 : Y:\1\DATA\011113\GB19133.D\FID1A.CH Vial: 20  
 Signal #2 : Y:\1\DATA\011113\GB19133.D\FID2B.CH  
 Acq On : 11 Jan 2013 10:35 pm Operator: StephK  
 Sample : D42556-2, 50X Inst : GC/MS Ins  
 Misc : GC3352,GGB1045,5.060,,100,5,1 Multiplr: 1.00  
 IntFile Signal #1: TVH1.E IntFile Signal #2: FB2.E  
 Quant Time: Jan 14 8:57 2013 Quant Results File: TB868GB868SOIL.RES

Quant Method : C:\MSDCHEM\1...\TB868GB868SOIL.M (Chemstation Integrator)  
 Title : 8015B/8021B TVH/BTEX  
 Last Update : Fri Jan 11 14:33:28 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: 23992846  
Conc: 0.35 mg/L m

#2 1,2,4-Trichlorobenzene

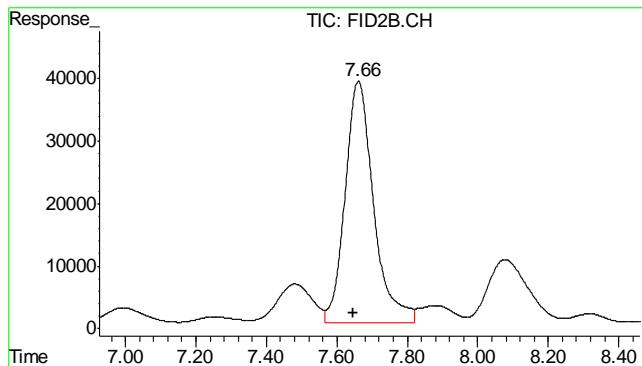
R.T.: 14.361 min  
Delta R.T.: 0.000 min  
Response: 2677351  
Conc: 85.45 % m

#4 Methyl-t-butyl-ether

R.T.: 0.000 min  
Exp R.T. : 2.241 min  
Response: 0  
Conc: N.D.

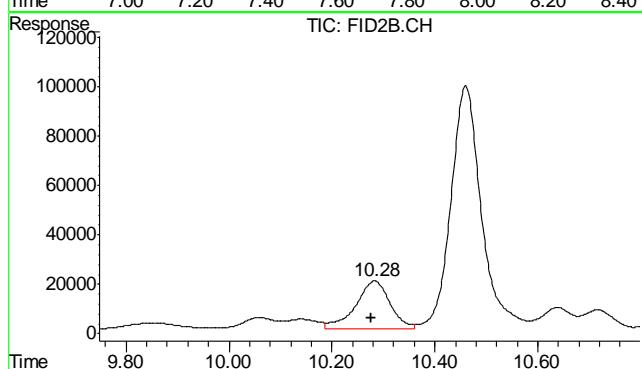
#5 Benzene

R.T.: 0.000 min  
Exp R.T. : 4.139 min  
Response: 0  
Conc: N.D.



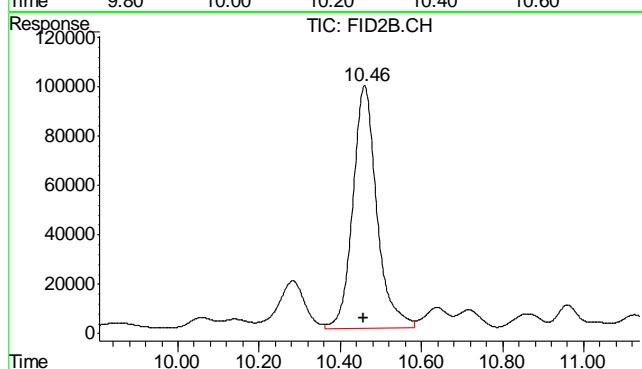
#6 Toluene

R.T.: 7.661 min  
Delta R.T.: 0.013 min  
Response: 2179890  
Conc: 5.50 ug/L



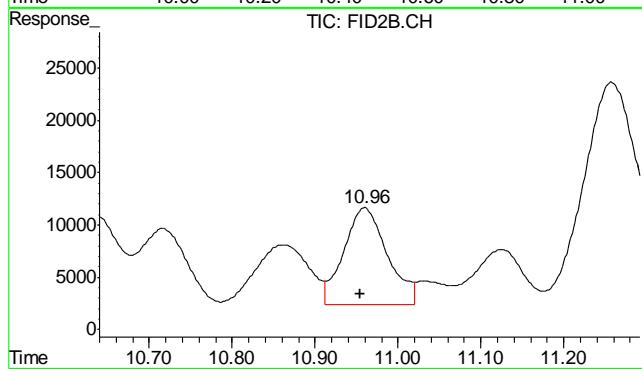
#7 Ethylbenzene

R.T.: 10.282 min  
Delta R.T.: 0.006 min  
Response: 888680  
Conc: 2.63 ug/L



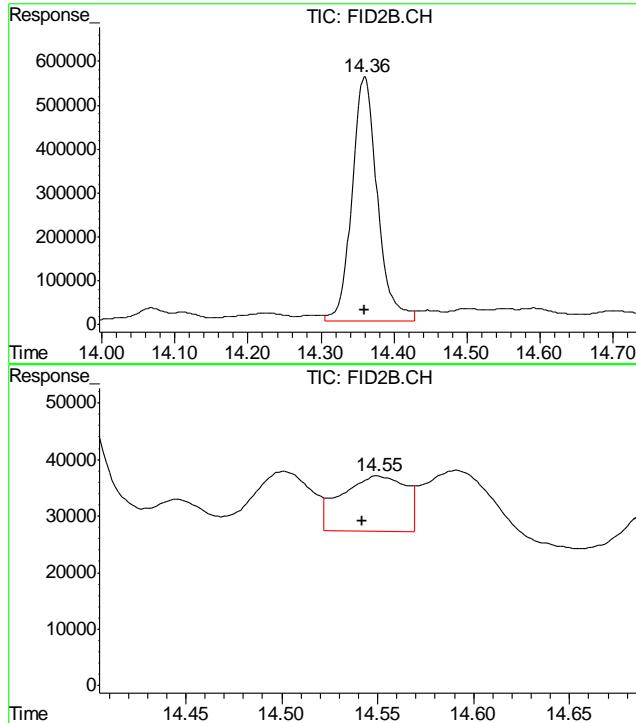
#8 m,p-Xylene

R.T.: 10.460 min  
Delta R.T.: 0.003 min  
Response: 4027842  
Conc: 10.66 ug/L



#9 o-Xylene

R.T.: 10.960 min  
Delta R.T.: 0.006 min  
Response: 334429  
Conc: 1.02 ug/L



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

R.T.: 14.359 min  
 Delta R.T.: 0.000 min  
 Response: 13368492  
 Conc: 82.25 % m

#11 Naphthalene

R.T.: 14.549 min  
 Delta R.T.: 0.008 min  
 Response: 226539  
 Conc: 1.15 ug/L m

11.1.2

11

**Manual Integrations  
APPROVED  
(compounds with "m" flag)**  
**Judy Nelson  
01/14/13 09:52**

## Quantitation Report (QT Reviewed)

Signal #1 : Y:\1\DATA\011113\GB19116.D\FID1A.CH Vial: 3  
 Signal #2 : Y:\1\DATA\011113\GB19116.D\FID2B.CH  
 Acq On : 11 Jan 2013 12:29 pm Operator: StephK  
 Sample : MB Inst : GC/MS Ins  
 Misc : GC3352,GGB1045,5.000,,100,5,1 Multiplr: 1.00  
 IntFile Signal #1: TVH1.E IntFile Signal #2: FB2.E  
 Quant Time: Jan 11 14:33:44 2013 Quant Results File: TB868GB868SOIL.RES

Quant Method : C:\MSDCHEM\1...\TB868GB868SOIL.M (Chemstation Integrator)  
 Title : 8015B/8021B TVH/BTEX  
 Last Update : Fri Jan 11 14:33:28 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
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**System Monitoring Compounds**

2) S	1,2,4-Trichlorobenzene	14.36	2820685	90.020 %	m
10) S	1,2,4-Trichlorobenzene (P)	14.36	14366912	88.397 %	m

**Target Compounds**

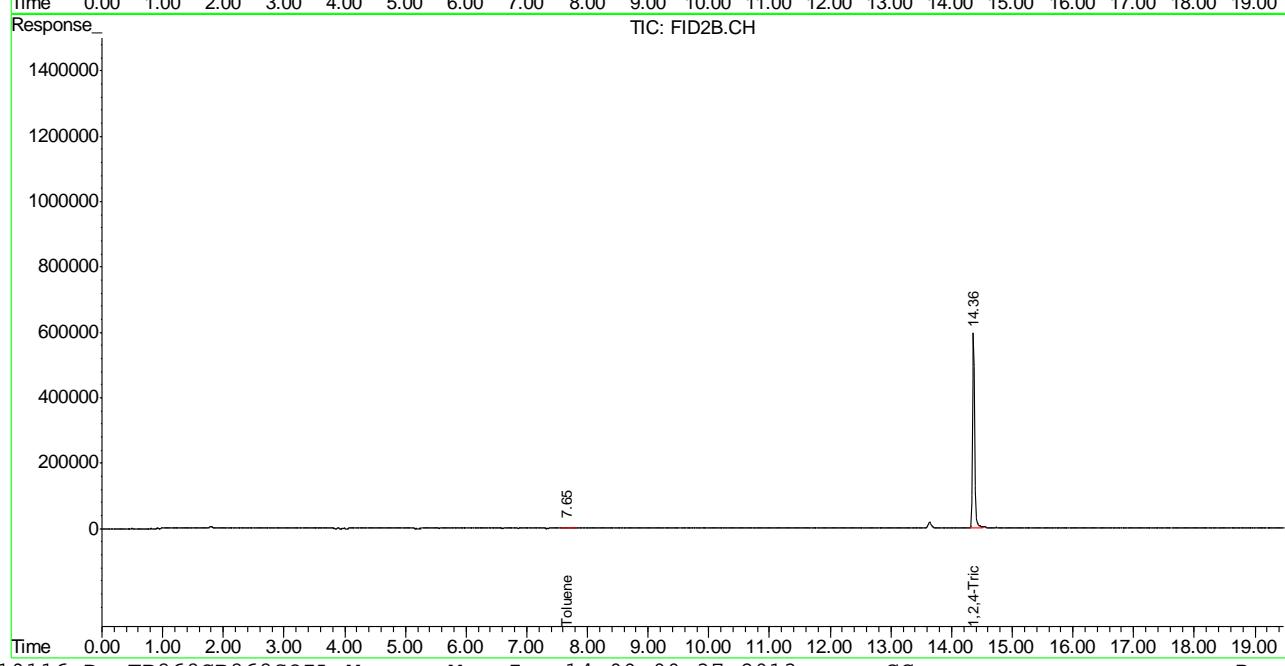
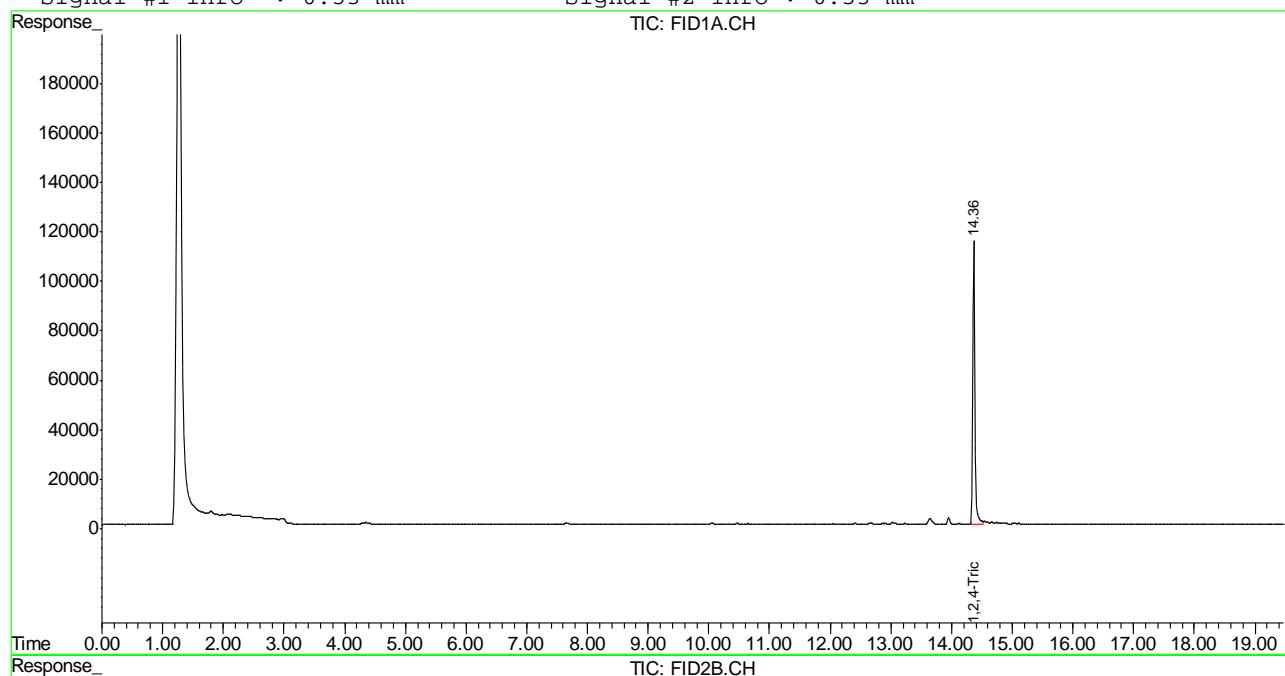
1) H	TVH-Gasoline	7.23	3021648	<MDL	mg/L
4) T	Methyl-t-butyl-ether	0.00	0	N.D.	ug/L d
5) T	Benzene	0.00	0	N.D.	ug/L d
6) T	Toluene	7.65	116878	0.295	ug/L
7) T	Ethylbenzene	0.00	0	N.D.	ug/L d
8) T	m,p-Xylene	0.00	0	N.D.	ug/L d
9) T	o-Xylene	0.00	0	N.D.	ug/L d
11) T	Naphthalene	14.55	16454	<MDL	ug/L m

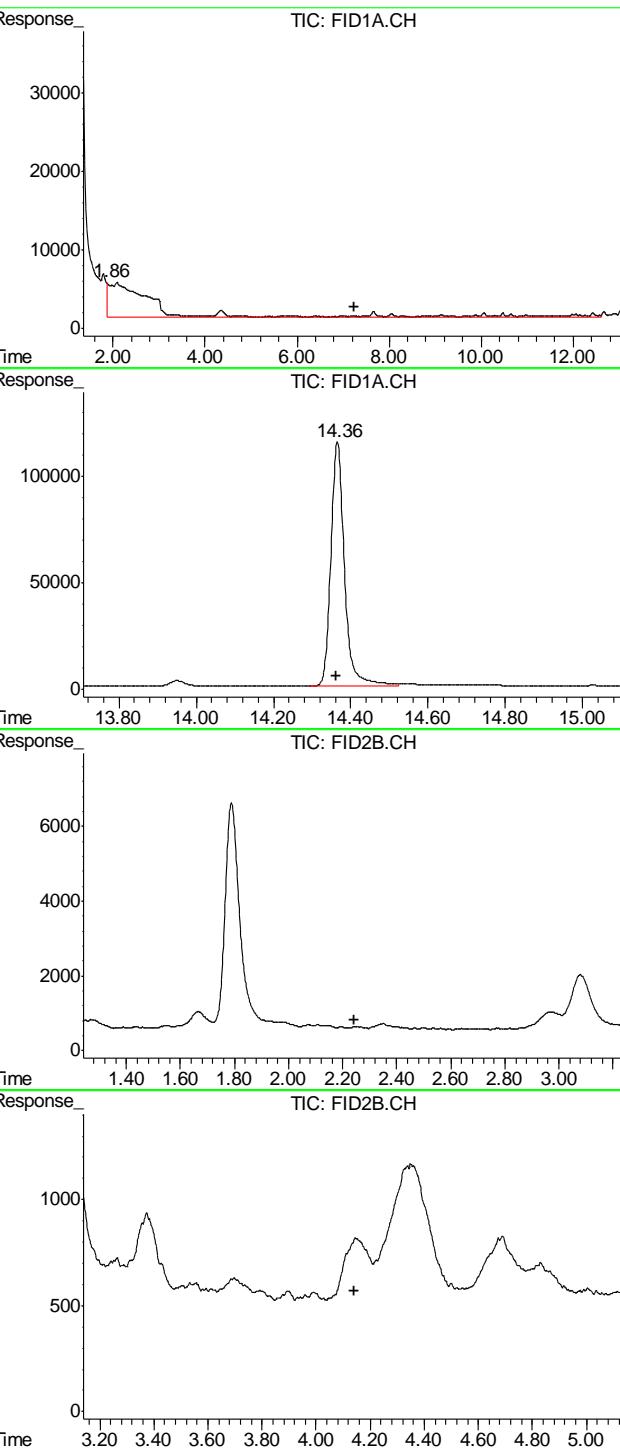
## Quantitation Report (QT Reviewed)

Signal #1 : Y:\1\DATA\011113\GB19116.D\FID1A.CH Vial: 3  
 Signal #2 : Y:\1\DATA\011113\GB19116.D\FID2B.CH  
 Acq On : 11 Jan 2013 12:29 pm Operator: StephK  
 Sample : MB Inst : GC/MS Ins  
 Misc : GC3352,GGB1045,5.000,,100,5,1 Multiplr: 1.00  
 IntFile Signal #1: TVH1.E IntFile Signal #2: FB2.E  
 Quant Time: Jan 11 14:33 2013 Quant Results File: TB868GB868SOIL.RES

Quant Method : C:\MSDCHEM\1...\TB868GB868SOIL.M (Chemstation Integrator)  
 Title : 8015B/8021B TVH/BTEX  
 Last Update : Fri Jan 11 14:33:28 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: 3021648  
 Conc: N.D.

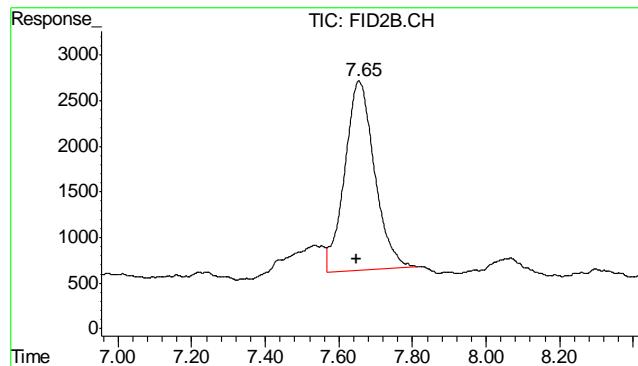
#2 1,2,4-Trichlorobenzene  
 R.T.: 14.365 min  
 Delta R.T.: 0.003 min  
 Response: 2820685  
 Conc: 90.02 % m

#4 Methyl-t-butyl-ether  
 R.T.: 0.000 min  
 Exp R.T. : 2.241 min  
 Response: 0  
 Conc: N.D.

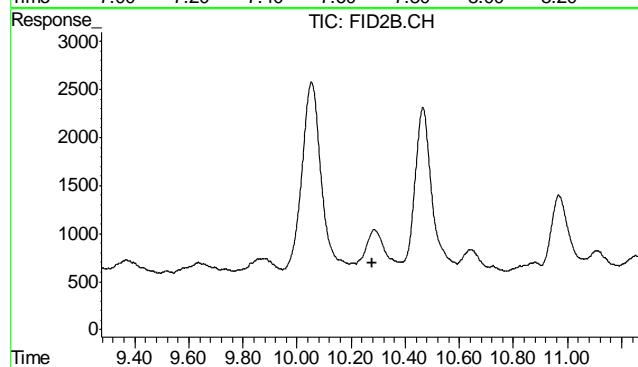
#5 Benzene  
 R.T.: 0.000 min  
 Exp R.T. : 4.139 min  
 Response: 0  
 Conc: N.D.

11.2.1

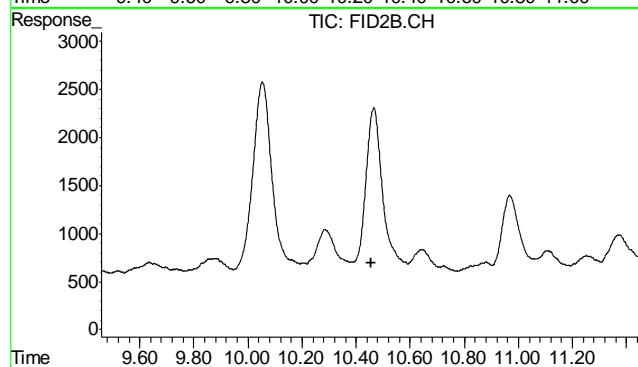
11



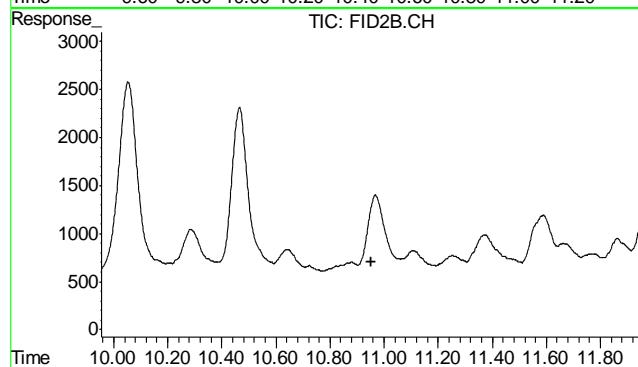
#6 Toluene  
R.T.: 7.655 min  
Delta R.T.: 0.006 min  
Response: 116878  
Conc: 0.29 ug/L



#7 Ethylbenzene  
R.T.: 0.000 min  
Exp R.T. : 10.276 min  
Response: 0  
Conc: N.D.



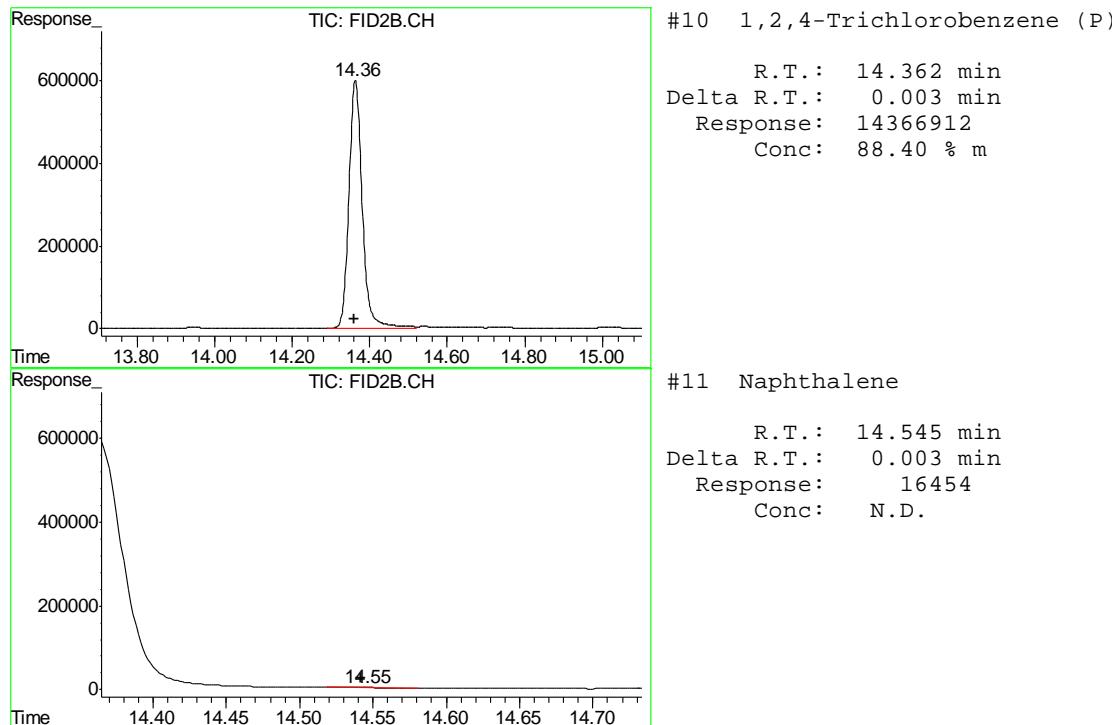
#8 m,p-Xylene  
R.T.: 0.000 min  
Exp R.T. : 10.457 min  
Response: 0  
Conc: N.D.



#9 o-Xylene  
R.T.: 0.000 min  
Exp R.T. : 10.954 min  
Response: 0  
Conc: N.D.

11.2.1

11



11.2.1

11



## GC Semi-volatiles

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### QC Data Summaries

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Includes the following where applicable:

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

**Method Blank Summary**

**Job Number:** D42556  
**Account:** XTOKWR XTO Energy  
**Project:** PCU 296-5A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
OP7232-MB	FD21147.D	1	01/15/13	AV	01/15/13	OP7232	GFD1064

The QC reported here applies to the following samples:

**Method:** SW846-8015B

D42556-1, D42556-2

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	79% 35-130%

## Blank Spike Summary

Page 1 of 1

Job Number: D42556

Account: XTOKWR XTO Energy

Project: PCU 296-5A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
OP7232-BS	FD21148.D	1	01/15/13	AV	01/15/13	OP7232	GFD1064

The QC reported here applies to the following samples:

Method: SW846-8015B

D42556-1, D42556-2

CAS No.	Compound	Spike mg/kg	BSP mg/kg	BSP %	Limits
	TPH-DRO (C10-C28)	667	631	95	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: D42556

Account: XTOKWR XTO Energy

Project: PCU 296-5A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
OP7232-MS	FD21149.D	1	01/15/13	AV	01/15/13	OP7232	GFD1064
OP7232-MSD	FD21150.D	1	01/15/13	AV	01/15/13	OP7232	GFD1064
D42562-1	FD21151.D	1	01/15/13	AV	01/15/13	OP7232	GFD1064

The QC reported here applies to the following samples:

Method: SW846-8015B

D42556-1, D42556-2

CAS No.	Compound	D42562-1		Spike mg/kg	MS mg/kg	MS %	MSD mg/kg	MSD %	RPD	Limits Rec/RPD
		mg/kg	Q							
	TPH-DRO (C10-C28)	313		876	1030	82	1030	82	0	20-168/30
CAS No.	Surrogate Recoveries	MS	MSD	D42562-1		Limits				
84-15-1	o-Terphenyl	91%	84%	79%						35-130%

\* = Outside of Control Limits.

12.3.1  
12



## GC Semi-volatiles

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Raw Data

---

## Quantitation Report (QT Reviewed)

Data File : C:\MSDCHEM\2\DATA\2013\JAN\FD011513\FD21152.D Vial: 8  
 Acq On : 1-15-2013 06:48:18 PM Operator: ashleyv  
 Sample : D42556-1 Inst : FID5  
 Misc : OP7232,GFD1064,30.10,,,1,1 Multiplr: 1.00  
 IntFile : autoint1.e  
 Quant Time: Jan 16 08:12:36 2013 Quant Results File: DRO-GFD1044F.RES

Quant Method : C:\MSDCHEM\2...\DRO-GFD1044F.M (Chemstation Integrator)  
 Title : 8015B TEH  
 Last Update : Tue Jan 15 13:27:58 2013  
 Response via : Initial Calibration  
 DataAcq Meth : DRO\_FR.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	8.81	80055375	1398.956 mg/L
<hr/>			
Target Compounds			
2) H TPH-DRO (C10-C32)	7.27	78643979	1918.275 mg/L
3) H TPH-DRO (C10-C28)	6.82	49423214	1210.711 mg/L

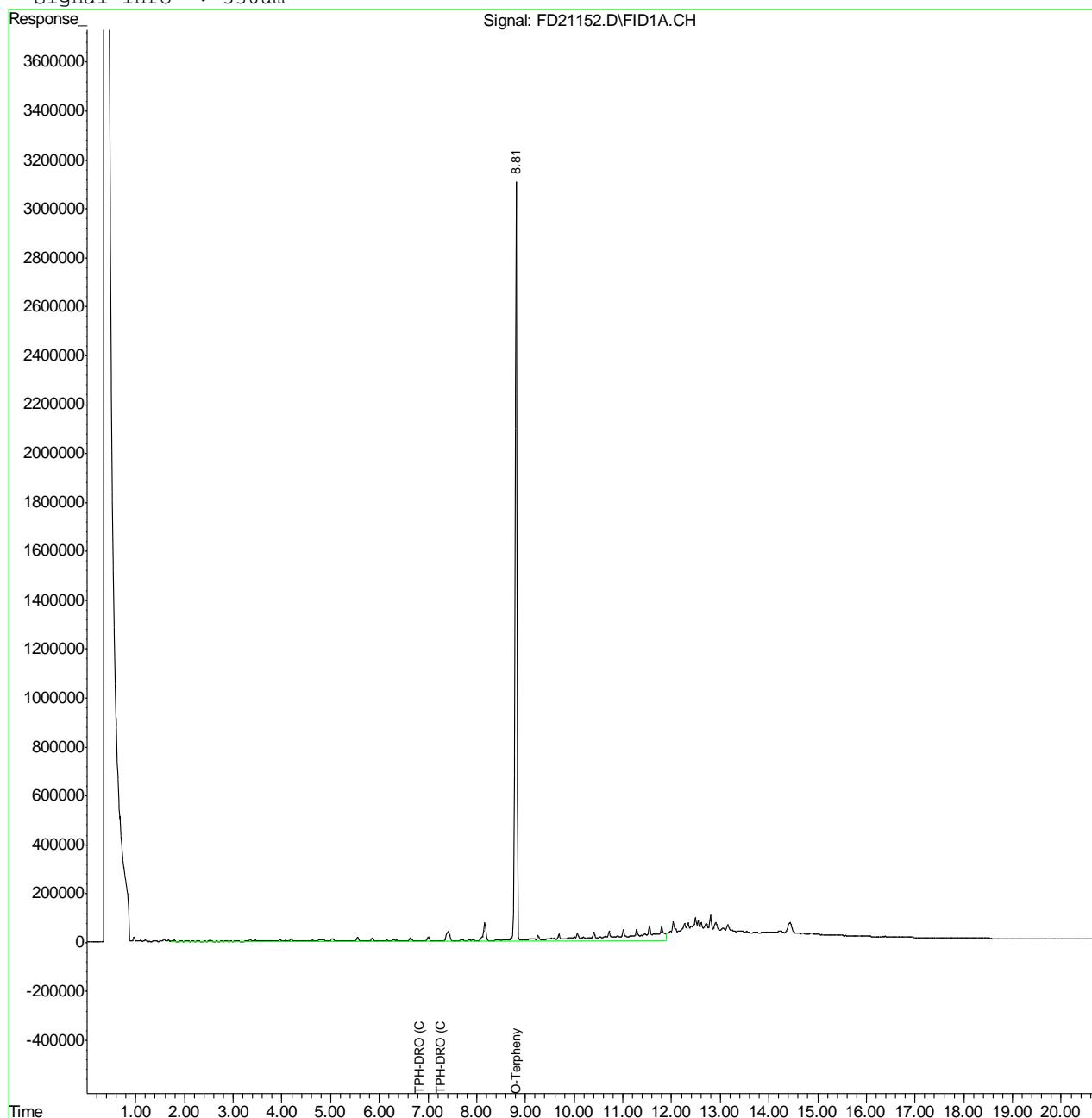
(f)=RT Delta > 1/2 Window (m)=manual int.  
 FD21152.D DRO-GFD1044F.M Wed Jan 16 08:34:02 2013 GC

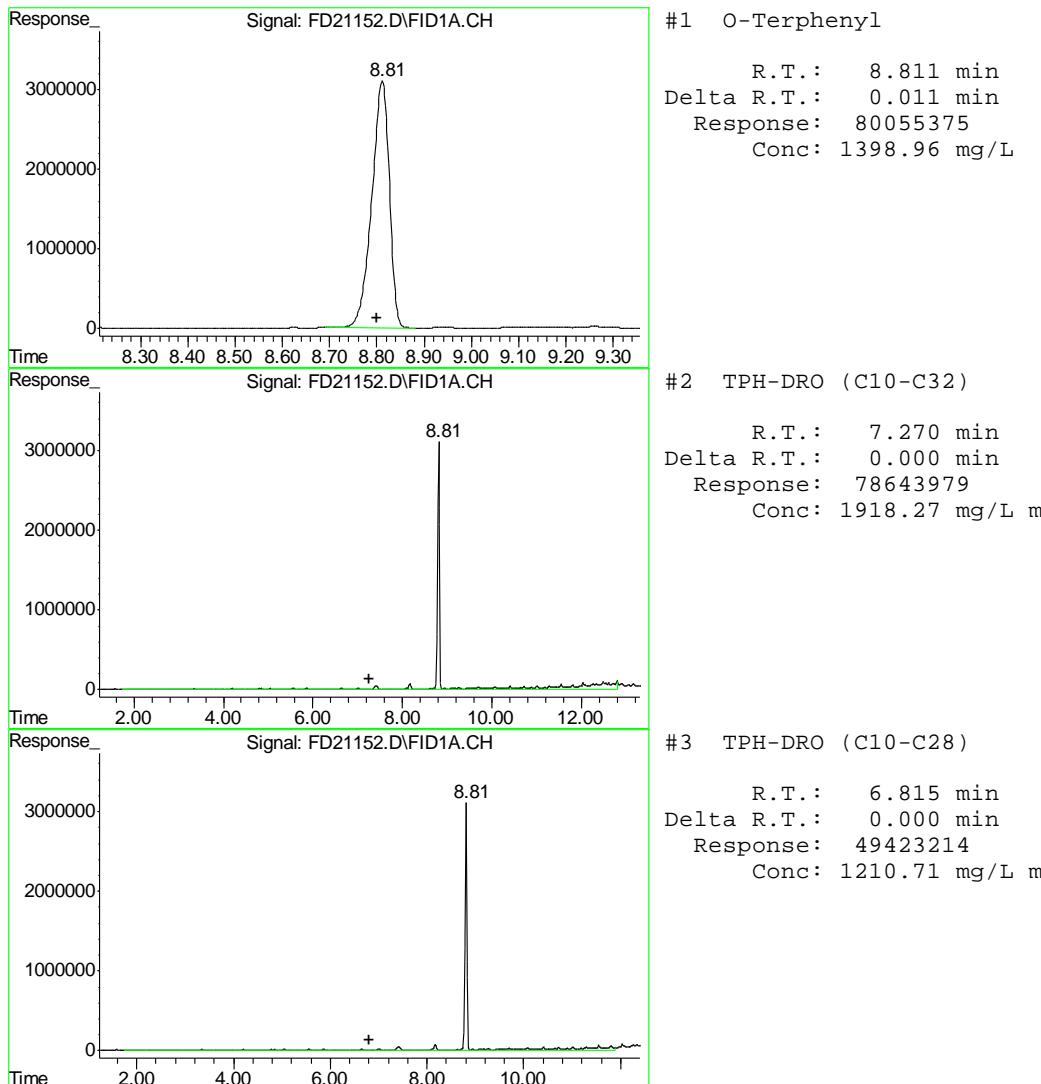
## Quantitation Report (QT Reviewed)

Data File : C:\MSDCHEM\2\DATA\2013\JAN\FD011513\FD21152.D Vial: 8  
 Acq On : 1-15-2013 06:48:18 PM Operator: ashleyv  
 Sample : D42556-1 Inst : FID5  
 Misc : OP7232,GFD1064,30.10,,,1,1 Multiplr: 1.00  
 IntFile : autoint1.e  
 Quant Time: Jan 16 8:12 2013 Quant Results File: DRO-GFD1044F.RES

Quant Method : C:\MSDCHEM\2...\DRO-GFD1044F.M (Chemstation Integrator)  
 Title : 8015B TEH  
 Last Update : Tue Jan 15 13:27:58 2013  
 Response via : Multiple Level Calibration  
 DataAcq Meth : DRO\_FR.M

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





## Quantitation Report (QT Reviewed)

Data File : C:\MSDCHEM\2\DATA\2013\JAN\FD011513\FD21153.D Vial: 9  
 Acq On : 1-15-2013 07:14:50 PM Operator: ashleyv  
 Sample : D42556-2 Inst : FID5  
 Misc : OP7232,GFD1064,30.01,,,1,1 Multiplr: 1.00  
 IntFile : autoint1.e  
 Quant Time: Jan 16 08:12:38 2013 Quant Results File: DRO-GFD1044F.RES

Quant Method : C:\MSDCHEM\2...\DRO-GFD1044F.M (Chemstation Integrator)  
 Title : 8015B TEH  
 Last Update : Tue Jan 15 13:27:58 2013  
 Response via : Initial Calibration  
 DataAcq Meth : DRO\_FR.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	8.81	92144707	1610.215 mg/L
<hr/>			
Target Compounds			
2) H TPH-DRO (C10-C32)	7.27	857459850	20915.058 mg/L
3) H TPH-DRO (C10-C28)	6.82	605699620	14837.711 mg/L

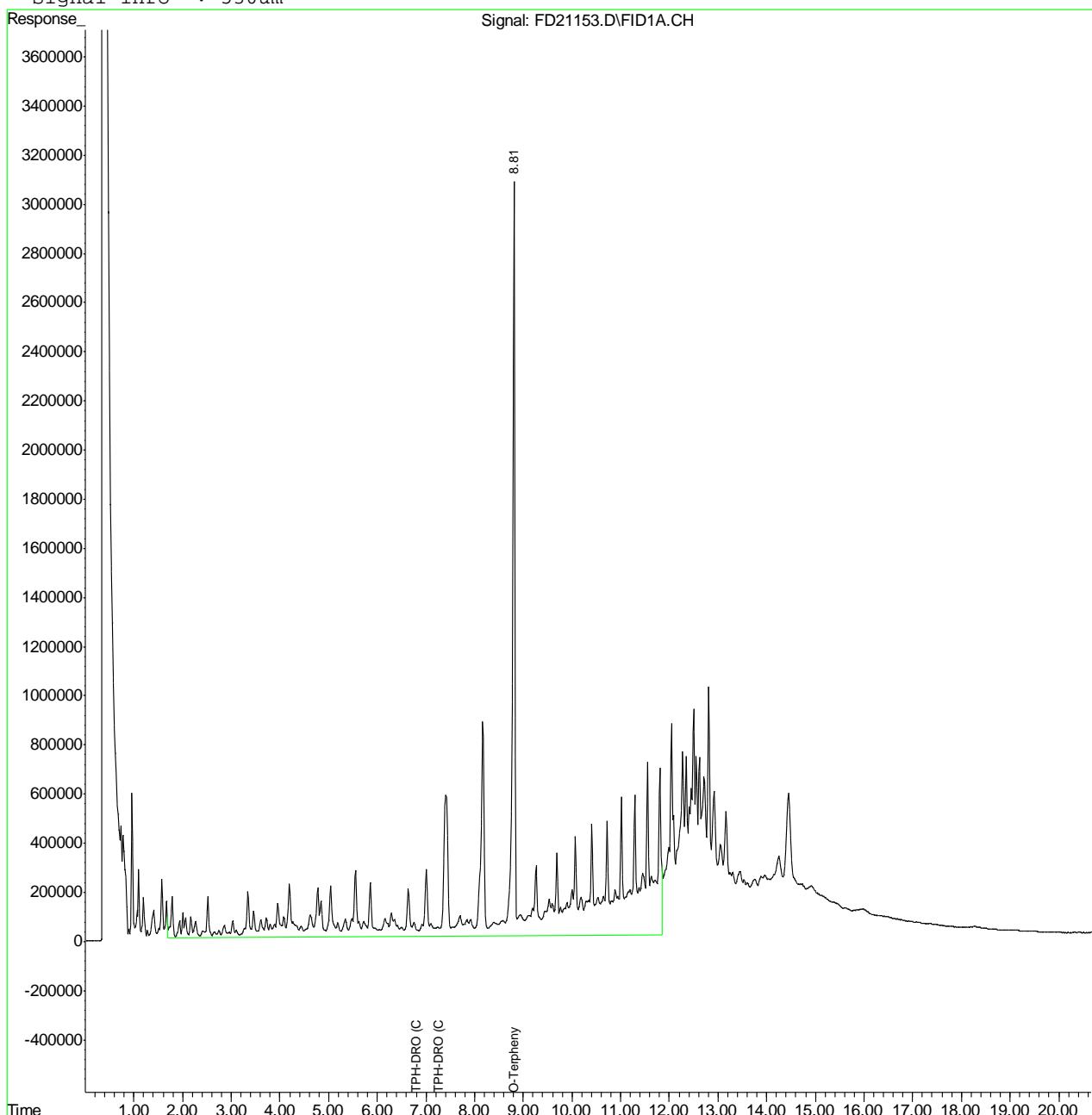
(f)=RT Delta > 1/2 Window (m)=manual int.  
 FD21153.D DRO-GFD1044F.M Wed Jan 16 08:34:03 2013 GC

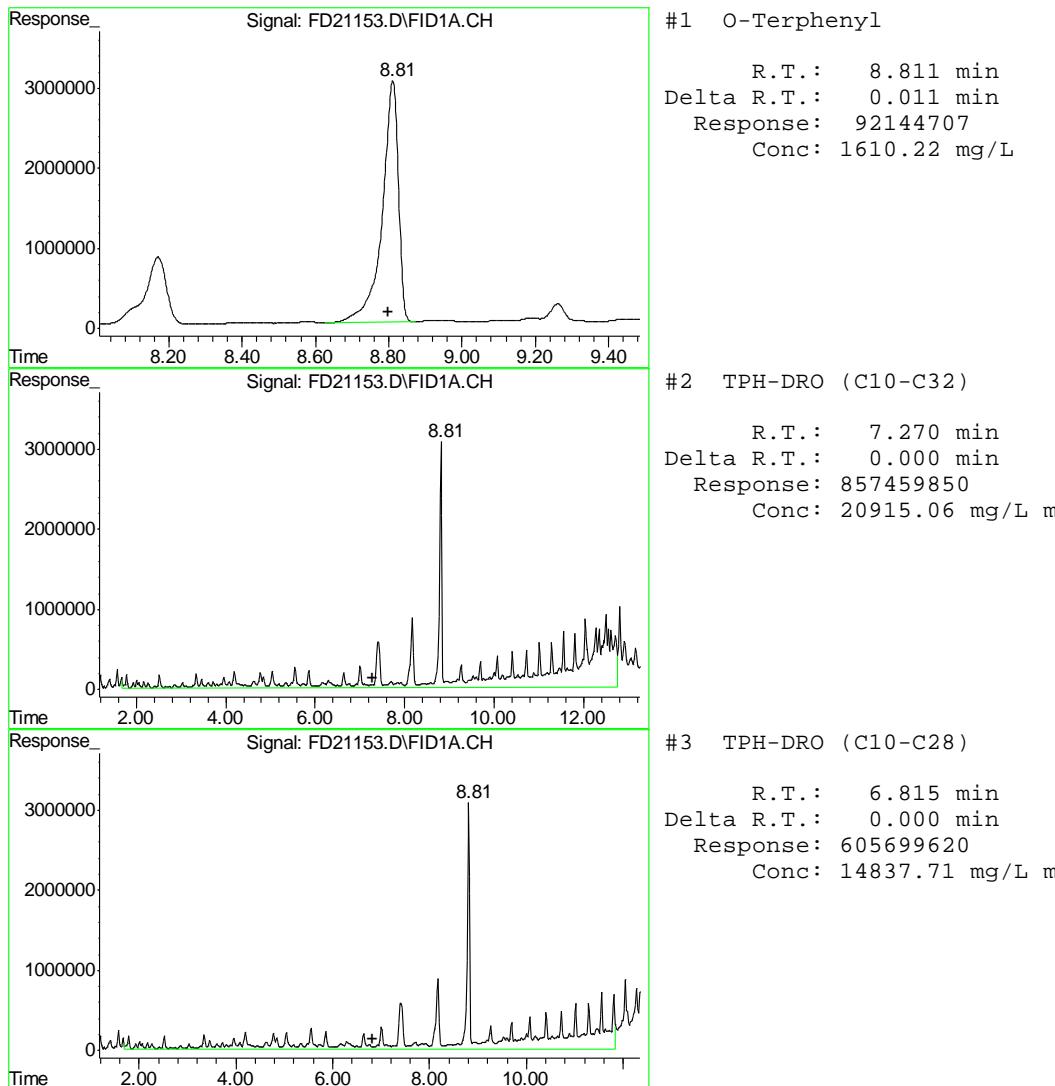
## Quantitation Report (QT Reviewed)

Data File : C:\MSDCHEM\2\DATA\2013\JAN\FD011513\FD21153.D Vial: 9  
 Acq On : 1-15-2013 07:14:50 PM Operator: ashleyv  
 Sample : D42556-2 Inst : FID5  
 Misc : OP7232,GFD1064,30.01,,,1,1 Multiplr: 1.00  
 IntFile : autoint1.e  
 Quant Time: Jan 16 8:23 2013 Quant Results File: DRO-GFD1044F.RES

Quant Method : C:\MSDCHEM\2...\DRO-GFD1044F.M (Chemstation Integrator)  
 Title : 8015B TEH  
 Last Update : Tue Jan 15 13:27:58 2013  
 Response via : Multiple Level Calibration  
 DataAcq Meth : DRO\_FR.M

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





## Quantitation Report (QT Reviewed)

Data File : C:\MSDCHEM\2\DATA\2013\JAN\FD011513\FD21147.D Vial: 3  
 Acq On : 1-15-2013 04:35:51 PM Operator: ashleyv  
 Sample : OP7232-MB Inst : FID5  
 Misc : OP7232,GFD1064,30.00,,,1,1 Multiplr: 1.00  
 IntFile : autoint1.e  
 Quant Time: Jan 16 08:12:28 2013 Quant Results File: DRO-GFD1044F.RES

Quant Method : C:\MSDCHEM\2...\DRO-GFD1044F.M (Chemstation Integrator)  
 Title : 8015B TEH  
 Last Update : Tue Jan 15 13:27:58 2013  
 Response via : Initial Calibration  
 DataAcq Meth : DRO\_FR.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	8.84	90567108	1582.647	mg/L
<hr/>				
Target Compounds				
2) H TPH-DRO (C10-C32)	7.27	2029145	49.495	mg/L
3) H TPH-DRO (C10-C28)	6.82	1512285	37.046	mg/L

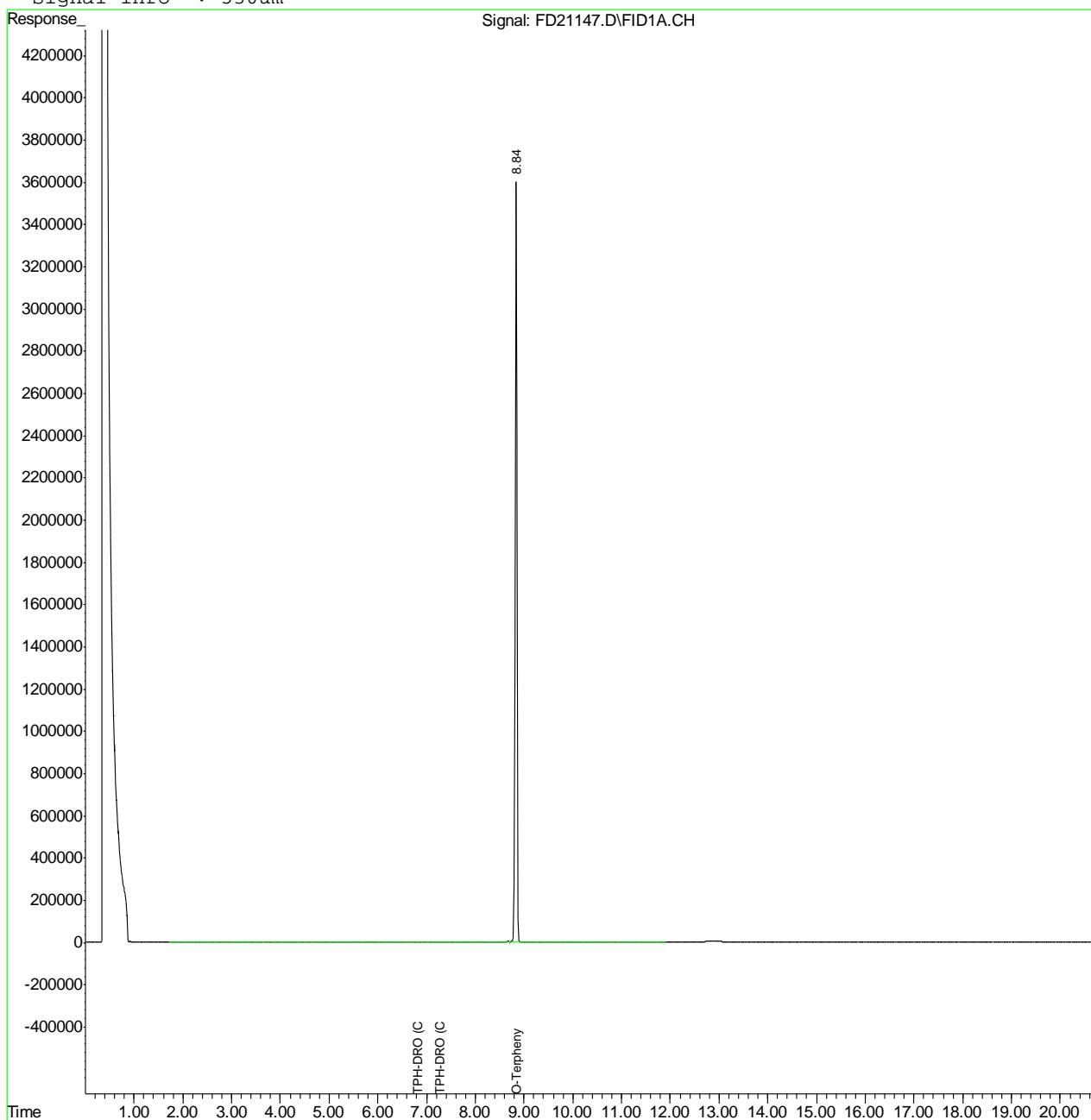
(f)=RT Delta > 1/2 Window (m)=manual int.  
 FD21147.D DRO-GFD1044F.M Wed Jan 16 08:33:57 2013 GC

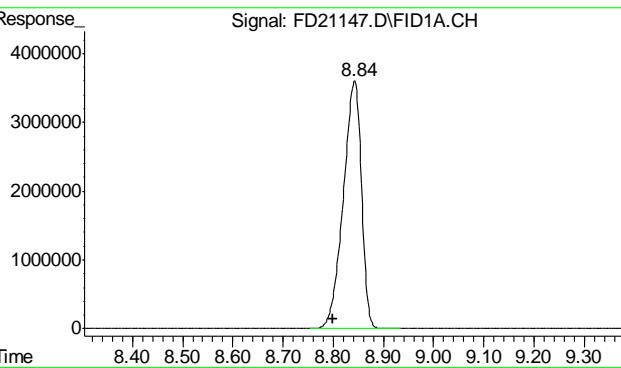
## Quantitation Report (QT Reviewed)

Data File : C:\MSDCHEM\2\DATA\2013\JAN\FD011513\FD21147.D Vial: 3  
 Acq On : 1-15-2013 04:35:51 PM Operator: ashleyv  
 Sample : OP7232-MB Inst : FID5  
 Misc : OP7232,GFD1064,30.00,,,1,1 Multiplr: 1.00  
 IntFile : autoint1.e  
 Quant Time: Jan 16 8:12 2013 Quant Results File: DRO-GFD1044F.RES

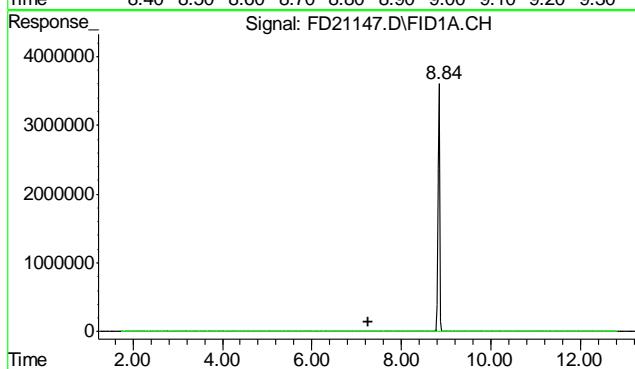
Quant Method : C:\MSDCHEM\2...\DRO-GFD1044F.M (Chemstation Integrator)  
 Title : 8015B TEH  
 Last Update : Tue Jan 15 13:27:58 2013  
 Response via : Multiple Level Calibration  
 DataAcq Meth : DRO\_FR.M

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

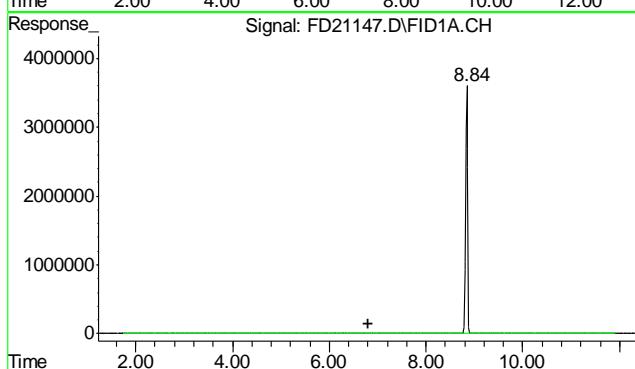




#1 O-Terphenyl  
R.T.: 8.842 min  
Delta R.T.: 0.042 min  
Response: 90567108  
Conc: 1582.65 mg/L



#2 TPH-DRO (C10-C32)  
R.T.: 7.270 min  
Delta R.T.: 0.000 min  
Response: 2029145  
Conc: 49.49 mg/L m



#3 TPH-DRO (C10-C28)  
R.T.: 6.815 min  
Delta R.T.: 0.000 min  
Response: 1512285  
Conc: 37.05 mg/L m



## Metals Analysis

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### QC Data Summaries

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Includes the following where applicable:

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

BLANK RESULTS SUMMARY  
Part 2 - Method Blanks

Login Number: D42556  
Account: XTOKRWR - XTO Energy  
Project: PCU 296-5A

QC Batch ID: MP9242  
Matrix Type: SOLID

Methods: SW846 6010C  
Units: mg/kg

Prep Date:

01/14/13

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	0.25	<1.0
Beryllium	1.0	.13	.15		
Boron	5.0	.1	.06		
Cadmium	1.0	.06	.036	0.020	<1.0
Calcium	40	.54	9		
Chromium	1.0	.03	.03	0.040	<1.0
Cobalt	0.50	.04	.07		
Copper	1.0	.12	.15	0.13	<1.0
Iron	7.0	.12	.87		
Lead	5.0	.19	.24	0.020	<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	-0.020	<3.0
Phosphorus	10	1.4	1.9		
Potassium	200	6.1	7		
Selenium	5.0	.48	.36	-0.46	<5.0
Silicon	5.0	.29	.37		
Silver	3.0	.04	.06	0.030	<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	0.090	<3.0

Associated samples MP9242: D42556-1, D42556-2

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

BLANK RESULTS SUMMARY  
Part 2 - Method Blanks

Login Number: D42556  
Account: XTOKRWR - XTO Energy  
Project: PCU 296-5A

QC Batch ID: MP9242  
Matrix Type: SOLID

Methods: SW846 6010C  
Units: mg/kg

Prep Date:

Metal

(anr) Analyte not requested

## MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D42556  
 Account: XTOKRWR - XTO Energy  
 Project: PCU 296-5A

QC Batch ID: MP9242  
 Matrix Type: SOLID

Methods: SW846 6010C  
 Units: mg/kg

Prep Date: 01/14/13

Metal	D42510-1 Original MS	Spikelot ICPALL2	% Rec	QC Limits
Aluminum				
Antimony				
Arsenic				
Barium	4420	4470	235	21.3 (a) 75-125
Beryllium				
Boron				
Cadmium	0.0	49.1	58.7	83.7 75-125
Calcium				
Chromium	28.6	73.0	58.7	83.2 75-125
Cobalt				
Copper	15.4	67.2	58.7	88.3 75-125
Iron				
Lead	12.0	107	117	81.0 75-125
Lithium				
Magnesium				
Manganese				
Molybdenum				
Nickel	19.2	77.2	58.7	98.9 75-125
Phosphorus				
Potassium				
Selenium	0.0	105	117	89.5 75-125
Silicon				
Silver	0.0	21.0	23.5	89.5 75-125
Sodium				
Strontium				
Thallium				
Tin				
Titanium				
Uranium				
Vanadium				
Zinc	37.6	82.7	58.7	76.9 75-125

Associated samples MP9242: D42556-1, D42556-2

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

MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D42556  
Account: XTOKRWR - XTO Energy  
Project: PCU 296-5A

QC Batch ID: MP9242  
Matrix Type: SOLID

Methods: SW846 6010C  
Units: mg/kg

Prep Date:

Metal

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

## MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D42556  
 Account: XTOKRWR - XTO Energy  
 Project: PCU 296-5A

QC Batch ID: MP9242  
 Matrix Type: SOLID

Methods: SW846 6010C  
 Units: mg/kg

Prep Date:

01/14/13

Metal	D42510-1 Original	MSD	Spikelot ICPALL2	% Rec	MSD RPD	QC Limit
Aluminum						
Antimony						
Arsenic						
Barium	4420	4930	232	219.5(a)	9.8	20
Beryllium						
Boron						
Cadmium	0.0	48.6	58.1	83.7	1.0	20
Calcium						
Chromium	28.6	74.5	58.1	86.6	2.0	20
Cobalt						
Copper	15.4	67.2	58.1	89.2	0.0	20
Iron						
Lead	12.0	105	116	80.1	1.9	20
Lithium						
Magnesium						
Manganese						
Molybdenum						
Nickel	19.2	77.0	58.1	99.5	0.3	20
Phosphorus						
Potassium						
Selenium	0.0	104	116	89.5	1.0	20
Silicon						
Silver	0.0	20.8	23.2	89.5	1.0	20
Sodium						
Strontium						
Thallium						
Tin						
Titanium						
Uranium						
Vanadium						
Zinc	37.6	81.9	58.1	76.3	1.0	20

Associated samples MP9242: D42556-1, D42556-2

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

MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D42556  
Account: XTOKRWR - XTO Energy  
Project: PCU 296-5A

QC Batch ID: MP9242  
Matrix Type: SOLID

Methods: SW846 6010C  
Units: mg/kg

Prep Date:

Metal

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

## SPIKE BLANK AND LAB CONTROL SAMPLE SUMMARY

Login Number: D42556  
 Account: XTOKRWR - XTO Energy  
 Project: PCU 296-5A

QC Batch ID: MP9242  
 Matrix Type: SOLID

Methods: SW846 6010C  
 Units: mg/kg

Prep Date: 01/14/13

Metal	BSP Result	Spikelot ICPALL2	% Rec	QC Limits
Aluminum				
Antimony				
Arsenic				
Barium	198	200	99.0	80-120
Beryllium				
Boron				
Cadmium	46.5	50	93.0	80-120
Calcium				
Chromium	49.9	50	99.8	80-120
Cobalt				
Copper	47.4	50	94.8	80-120
Iron				
Lead	96.0	100	96.0	80-120
Lithium				
Magnesium				
Manganese				
Molybdenum				
Nickel	47.5	50	95.0	80-120
Phosphorus				
Potassium				
Selenium	97.5	100	97.5	80-120
Silicon				
Silver	19.8	20	99.0	80-120
Sodium				
Strontium				
Thallium				
Tin				
Titanium				
Uranium				
Vanadium				
Zinc	49.2	50	98.4	80-120

Associated samples MP9242: D42556-1, D42556-2

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

SPIKE BLANK AND LAB CONTROL SAMPLE SUMMARY

Login Number: D42556  
Account: XTOKRWR - XTO Energy  
Project: PCU 296-5A

QC Batch ID: MP9242  
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: D42556  
 Account: XTOKRWR - XTO Energy  
 Project: PCU 296-5A

QC Batch ID: MP9242  
 Matrix Type: SOLID

Methods: SW846 6010C  
 Units: ug/l

Prep Date:

01/14/13

Metal	D42510-1 Original	SDL 1:5	%DIF	QC Limits
Aluminum				
Antimony				
Arsenic				
Barium	32800	36700	2.5	0-10
Beryllium				
Boron				
Cadmium	0.00	0.00	NC	0-10
Calcium				
Chromium	206	243	17.8*(a)	0-10
Cobalt				
Copper	121	142	8.3	0-10
Iron				
Lead	119	122	19.0*(a)	0-10
Lithium				
Magnesium				
Manganese				
Molybdenum				
Nickel	135	155	5.2	0-10
Phosphorus				
Potassium				
Selenium	0.00	0.00	NC	0-10
Silicon				
Silver	0.00	0.00	NC	0-10
Sodium				
Strontium				
Thallium				
Tin				
Titanium				
Uranium				
Vanadium				
Zinc	320	412	28.5*(a)	0-10

Associated samples MP9242: D42556-1, D42556-2

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

SERIAL DILUTION RESULTS SUMMARY

Login Number: D42556  
Account: XTOKRWR - XTO Energy  
Project: PCU 296-5A

QC Batch ID: MP9242  
Matrix Type: SOLID

Methods: SW846 6010C  
Units: ug/l

Prep Date:

Metal

(anr) Analyte not requested  
(a) Serial dilution indicates possible matrix interference.

BLANK RESULTS SUMMARY  
Part 2 - Method Blanks

Login Number: D42556  
Account: XTOKWR - XTO Energy  
Project: PCU 296-5A

QC Batch ID: MP9243  
Matrix Type: SOLID

Methods: SW846 6020A  
Units: mg/kg

Prep Date:

01/14/13

Metal	RL	IDL	MDL	MB raw	final
Aluminum	25	.22	.31		
Antimony	0.20	.0018	.0075		
Arsenic	0.10	.006	.06	0.013	<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 MP9243: D42556-1, D42556-2

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: D42556  
 Account: XTOKWR - XTO Energy  
 Project: PCU 296-5A

QC Batch ID: MP9243  
 Matrix Type: SOLID

Methods: SW846 6020A  
 Units: mg/kg

Prep Date:

01/14/13

Metal	D42510-1 Original MS	Spikelot ICPALL2	% Rec	QC Limits
Aluminum				
Antimony				
Arsenic	11.1	116	117	89.4    75-125
Barium	anr			
Beryllium				
Boron				
Cadmium	anr			
Calcium				
Chromium	anr			
Cobalt				
Copper				
Iron				
Lead	anr			
Magnesium				
Manganese				
Molybdenum				
Nickel				
Phosphorus				
Potassium				
Selenium	anr			
Silver	anr			
Sodium				
Strontium				
Thallium				
Tin				
Titanium				
Uranium				
Vanadium				
Zinc				

Associated samples MP9243: D42556-1, D42556-2

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: D42556  
 Account: XTOKWR - XTO Energy  
 Project: PCU 296-5A

QC Batch ID: MP9243  
 Matrix Type: SOLID

Methods: SW846 6020A  
 Units: mg/kg

Prep Date: 01/14/13

Metal	D42510-1 Original	MSD	Spikelot ICPALL2	% Rec	MSD RPD	QC Limit
Aluminum						
Antimony						
Arsenic	11.1	123	116	96.3	5.9	20
Barium						
Beryllium						
Boron						
Cadmium		anr				
Calcium						
Chromium		anr				
Cobalt						
Copper						
Iron						
Lead		anr				
Magnesium						
Manganese						
Molybdenum						
Nickel						
Phosphorus						
Potassium						
Selenium		anr				
Silver		anr				
Sodium						
Strontium						
Thallium						
Tin						
Titanium						
Uranium						
Vanadium						
Zinc						

Associated samples MP9243: D42556-1, D42556-2

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: D42556  
 Account: XTOKRWR - XTO Energy  
 Project: PCU 296-5A

QC Batch ID: MP9243  
 Matrix Type: SOLID

Methods: SW846 6020A  
 Units: mg/kg

Prep Date: 01/14/13

Metal	BSP Result	Spikelot ICPALL2	% Rec	QC Limits
Aluminum				
Antimony				
Arsenic	101	100	101.0	80-120
Barium	anr			
Beryllium				
Boron				
Cadmium	anr			
Calcium				
Chromium	anr			
Cobalt				
Copper				
Iron				
Lead	anr			
Magnesium				
Manganese				
Molybdenum				
Nickel				
Phosphorus				
Potassium				
Selenium	anr			
Silver	anr			
Sodium				
Strontium				
Thallium				
Tin				
Titanium				
Uranium				
Vanadium				
Zinc				

Associated samples MP9243: D42556-1, D42556-2

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

14.2.3  
**14**

## SERIAL DILUTION RESULTS SUMMARY

Login Number: D42556  
 Account: XTOKRWR - XTO Energy  
 Project: PCU 296-5A

QC Batch ID: MP9243  
 Matrix Type: SOLID

Methods: SW846 6020A  
 Units: ug/l

Prep Date: 01/14/13

Metal	D42510-1 Original	SDL 5:25	%DIF	QC Limits
Aluminum				
Antimony				
Arsenic	94.9	94.9	0.0	0-10
Barium	anr			
Beryllium				
Boron				
Cadmium	anr			
Calcium				
Chromium	anr			
Cobalt				
Copper				
Iron				
Lead	anr			
Magnesium				
Manganese				
Molybdenum				
Nickel				
Phosphorus				
Potassium				
Selenium	anr			
Silver	anr			
Sodium				
Strontium				
Thallium				
Tin				
Titanium				
Uranium				
Vanadium				
Zinc				

Associated samples MP9243: D42556-1, D42556-2

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: D42556  
Account: XTOKRWR - XTO Energy  
Project: PCU 296-5A

QC Batch ID: MP9244  
Matrix Type: SOLID

Methods: SW846 7471B  
Units: mg/kg

Prep Date: 01/15/13

Metal	RL	IDL	MDL	MB raw	final
Mercury	0.10	.0011	.0009	0.0018	<0.10

Associated samples MP9244: D42556-1, D42556-2

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

14.3.1  
**14**

MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D42556  
Account: XTOKRWR - XTO Energy  
Project: PCU 296-5A

QC Batch ID: MP9244  
Matrix Type: SOLID

Methods: SW846 7471B  
Units: mg/kg

Prep Date: 01/15/13

Metal	D42445-1 Original MS	Spikelot HGWSR1	QC % Rec	QC Limits
Mercury	0.054	0.45	0.393	100.7 75-125

Associated samples MP9244: D42556-1, D42556-2

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: D42556  
Account: XTOKRWR - XTO Energy  
Project: PCU 296-5A

QC Batch ID: MP9244  
Matrix Type: SOLID

Methods: SW846 7471B  
Units: mg/kg

Prep Date:

01/15/13

Metal	D42445-1 Original	MSD	Spikelot HGWSR1	MSD % Rec	QC RPD	QC Limit
Mercury	0.054	0.45	0.393	100.7	0.0	20

Associated samples MP9244: D42556-1, D42556-2

Results &lt; 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: D42556  
Account: XTOKRWR - XTO Energy  
Project: PCU 296-5A

QC Batch ID: MP9244  
Matrix Type: SOLID

Methods: SW846 7471B  
Units: mg/kg

Prep Date: 01/15/13

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

Associated samples MP9244: D42556-1, D42556-2

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: D42556  
Account: XTOKRWR - XTO Energy  
Project: PCU 296-5A

QC Batch ID: MP9251  
Matrix Type: AQUEOUS

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

Prep Date:

01/15/13

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	2.5	<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	24.0	<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	27.5	<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 MP9251: D42556-1A, D42556-2A

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

BLANK RESULTS SUMMARY  
Part 2 - Method Blanks

Login Number: D42556  
Account: XTOKRWR - XTO Energy  
Project: PCU 296-5A

QC Batch ID: MP9251  
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: D42556  
 Account: XTOKWR - XTO Energy  
 Project: PCU 296-5A

QC Batch ID: MP9251  
 Matrix Type: AQUEOUS

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

Prep Date:

01/15/13

Metal	D42556-1A Original MS	Spikelot ICPALL2	% Rec	QC Limits
Aluminum				
Antimony				
Arsenic				
Barium				
Beryllium				
Boron				
Cadmium				
Calcium	92300	231000	125000	111.0
Chromium				
Cobalt				
Copper				
Iron				
Lead				
Lithium				
Magnesium	26000	155000	125000	103.2
Manganese				
Molybdenum				
Nickel				
Phosphorus				
Potassium				
Selenium				
Silicon				
Silver				
Sodium	107000	239000	125000	105.6
Strontium				
Thallium				
Tin				
Titanium				
Uranium				
Vanadium				
Zinc				

Associated samples MP9251: D42556-1A, D42556-2A

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

MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D42556  
Account: XTOKRWR - XTO Energy  
Project: PCU 296-5A

QC Batch ID: MP9251  
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: D42556  
 Account: XTOKRWR - XTO Energy  
 Project: PCU 296-5A

QC Batch ID: MP9251  
 Matrix Type: AQUEOUS

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

Prep Date: 01/15/13

Metal	D42556-1A Original MSD	Spikelot ICPALL2	MSD % Rec	MSD RPD	QC Limit
Aluminum					
Antimony					
Arsenic					
Barium					
Beryllium					
Boron					
Cadmium					
Calcium	92300	236000	125000	115.0	2.1
Chromium					
Cobalt					
Copper					
Iron					
Lead					
Lithium					
Magnesium	26000	153000	125000	101.6	1.3
Manganese					
Molybdenum					
Nickel					
Phosphorus					
Potassium					
Selenium					
Silicon					
Silver					
Sodium	107000	235000	125000	102.4	1.7
Strontium					
Thallium					
Tin					
Titanium					
Uranium					
Vanadium					
Zinc					

Associated samples MP9251: D42556-1A, D42556-2A

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

MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D42556  
Account: XTOKRWR - XTO Energy  
Project: PCU 296-5A

QC Batch ID: MP9251  
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: D42556  
 Account: XTOKRWR - XTO Energy  
 Project: PCU 296-5A

QC Batch ID: MP9251  
 Matrix Type: AQUEOUS

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

Prep Date: 01/15/13

Metal	BSP Result	Spikelot ICPALL2	% Rec	QC Limits
Aluminum				
Antimony				
Arsenic				
Barium				
Beryllium				
Boron				
Cadmium				
Calcium	137000	125000	109.6	80-120
Chromium				
Cobalt				
Copper				
Iron				
Lead				
Lithium				
Magnesium	123000	125000	98.4	80-120
Manganese				
Molybdenum				
Nickel				
Phosphorus				
Potassium				
Selenium				
Silicon				
Silver				
Sodium	125000	125000	100.0	80-120
Strontium				
Thallium				
Tin				
Titanium				
Uranium				
Vanadium				
Zinc				

Associated samples MP9251: D42556-1A, D42556-2A

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

SPIKE BLANK AND LAB CONTROL SAMPLE SUMMARY

Login Number: D42556  
Account: XTOKRWR - XTO Energy  
Project: PCU 296-5A

QC Batch ID: MP9251  
Matrix Type: AQUEOUS

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

Prep Date:

Metal

(anr) Analyte not requested

## SERIAL DILUTION RESULTS SUMMARY

Login Number: D42556  
 Account: XTOKRWR - XTO Energy  
 Project: PCU 296-5A

QC Batch ID: MP9251  
 Matrix Type: AQUEOUS

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

Prep Date: 01/15/13

Metal	D42556-1A Original	SDL 1:5	%DIF	QC Limits
Aluminum				
Antimony				
Arsenic				
Barium				
Beryllium				
Boron				
Cadmium				
Calcium	18500	18600	0.9	0-10
Chromium				
Cobalt				
Copper				
Iron				
Lead				
Lithium				
Magnesium	5190	5270	1.5	0-10
Manganese				
Molybdenum				
Nickel				
Phosphorus				
Potassium				
Selenium				
Silicon				
Silver				
Sodium	21500	21600	0.8	0-10
Strontium				
Thallium				
Tin				
Titanium				
Uranium				
Vanadium				
Zinc				

Associated samples MP9251: D42556-1A, D42556-2A

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

SERIAL DILUTION RESULTS SUMMARY

Login Number: D42556  
Account: XTOKRWR - XTO Energy  
Project: PCU 296-5A

QC Batch ID: MP9251  
Matrix Type: AQUEOUS

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

Prep Date:

Metal

(anr) Analyte not requested



## General Chemistry

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### QC Data Summaries

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Includes the following where applicable:

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

METHOD BLANK AND SPIKE RESULTS SUMMARY  
GENERAL CHEMISTRY

Login Number: D42556  
Account: XTOKWR - XTO Energy  
Project: PCU 296-5A

Analyte	Batch ID	RL	MB Result	Units	Spike Amount	BSP Result	BSP %Recov	QC Limits
Chromium, Hexavalent	GP9086/GN18425	1.0	0.0	mg/kg	92.9	91.5	98.5	80-120%
Specific Conductivity	GP9098/GN18435	1.0	<1.0	umhos/cm	9992	10500	105.2	90-110%
pH	GN18424			su	8.00	7.97	99.8	99.3-100.7%
pH	GN18426			su	8.00	8.00	100.0	99.3-100.7%

Associated Samples:

Batch GP9086: D42556-1, D42556-2

Batch GP9098: D42556-1, D42556-2

Batch GN18424: D42556-1

Batch GN18426: D42556-2

(\*) Outside of QC limits

DUPLICATE RESULTS SUMMARY  
GENERAL CHEMISTRY

Login Number: D42556  
Account: XTOKWR - XTO Energy  
Project: PCU 296-5A

Analyte	Batch ID	QC Sample	Units	Original Result	DUP Result	RPD	QC Limits
Chromium, Hexavalent Redox Potential Vs H2	GP9086/GN18425 GN18422	D42556-1 D42337-20	mg/kg mv	0.0 230	0.0 227	0.0 1.3	0-20% 0-20%

Associated Samples:

Batch GP9086: D42556-1, D42556-2  
Batch GN18422: D42556-1, D42556-2

(\*) Outside of QC limits

MATRIX SPIKE RESULTS SUMMARY  
GENERAL CHEMISTRY

Login Number: D42556  
Account: XTOKWR - XTO Energy  
Project: PCU 296-5A

Analyte	Batch ID	QC Sample	Units	Original Result	Spike Amount	MS Result	%Rec	QC Limits
Chromium, Hexavalent	GP9086/GN18425	D42556-1	mg/kg	0.0	40.0	35.0	87.5	75-125%

Associated Samples:

Batch GP9086: D42556-1, D42556-2

(\*) Outside of QC limits

(N) Matrix Spike Rec. outside of QC limits

MATRIX SPIKE DUPLICATE RESULTS SUMMARY  
GENERAL CHEMISTRY

Login Number: D42556  
Account: XTOKWR - XTO Energy  
Project: PCU 296-5A

Analyte	Batch ID	QC Sample	Units	Original Result	Spike Amount	MSD Result	RPD	QC Limit
Chromium, Hexavalent	GP9086/GN18425	D42556-1	mg/kg	0.0	40.0	33.9	3.1	20%

Associated Samples:

Batch GP9086: D42556-1, D42556-2

(\*) Outside of QC limits

(N) Matrix Spike Rec. outside of QC limits