

**Technical Report for**

**XTO Energy**

**FRU 297-8B**

**1106-06**

**Accutest Job Number: D35710**

**Sampling Date: 06/20/12**

**Report to:**

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

**Total number of pages in report: 150**



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.



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

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

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

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Test results relate only to samples analyzed.

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## Sample Summary

XTO Energy

Job No: D35710

FRU 297-8B

Project No: 1106-06

Sample Number	Collected		Received	Matrix		Client Sample ID
	Date	Time By		Code	Type	
D35710-1	06/20/12	11:45 DS	06/21/12	SO	Soil	CUT 1 SUBLINER
D35710-1A	06/20/12	11:45 DS	06/21/12	SO	Soil	CUT 1 SUBLINER

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



## CASE NARRATIVE / CONFORMANCE SUMMARY

**Client:** XTO Energy

**Job No** D35710

**Site:** FRU 297-8B

**Report Date** 6/28/2012 8:51:31 AM

On 06/21/2012, 1 sample(s), 0 Trip Blank(s), and 0 Field Blank(s) were received at Accutest Mountain States (AMS) at a temperature of 4.0 °C. The samples were intact and properly preserved, unless noted below. An AMS Job Number of D35710 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> V5V1355
------------------	--------------------------

- All samples were analyzed within the recommended method holding time.
- All method blanks for this batch meet method specific criteria.
- Sample(s) D35708-1MS, D35708-1MSD were used as the QC samples indicated.

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

<b>Matrix</b> SO	<b>Batch ID:</b> OP6113
------------------	-------------------------

- All samples were extracted and analyzed within the recommended method holding time.
- Sample(s) D35708-1MS, D35708-1MSD were used as the QC samples indicated.
- All method blanks for this batch meet method specific criteria.

### Volatiles by GC By Method SW846 8015B

<b>Matrix</b> SO	<b>Batch ID:</b> GGB910
------------------	-------------------------

- All samples were analyzed within the recommended method holding time.
- All method blanks for this batch meet method specific criteria.
- Sample(s) D35708-1MS, D35708-1MSD were used as the QC samples indicated.

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

<b>Matrix</b> SO	<b>Batch ID:</b> OP6112
------------------	-------------------------

- All samples were extracted and analyzed within the recommended method holding time.
- Sample(s) D35708-1MS, D35708-1MSD 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:** MP7743

- All samples were digested and analyzed within the recommended method holding time.
- All method blanks for this batch meet method specific criteria.
- Sample(s) D35708-1AMS, D35708-1AMSD, D35708-1ASDL were used as the QC samples for the metals analysis.
- The matrix spike (MS) recovery(s) of Sodium are outside control limits. Spike amount low relative to the sample amount. Refer to lab control or spike blank for recovery information.

**Matrix** SO **Batch ID:** MP7737

- All samples were digested and analyzed within the recommended method holding time.
- All method blanks for this batch meet method specific criteria.
- Sample(s) D35708-1MS, D35708-1MSD, D35708-1SDL were used as the QC samples for the metals analysis.
- The matrix spike (MS) recovery(s) of Barium are outside control limits. Spike recovery indicates possible matrix interference.
- The serial dilution RPD(s) for Cadmium, Selenium, Barium, Chromium, Nickel, Zinc are outside control limits for sample MP7737-SD1. Percent difference acceptable due to low initial sample concentration (< 50 times IDL).
- MP7737-SD1 for Barium, Chromium, Nickel, and Zinc: Serial dilution indicates possible matrix interference.

## Metals By Method SW846 6020A

**Matrix** SO **Batch ID:** MP7738

- All samples were digested and analyzed within the recommended method holding time.
- All method blanks for this batch meet method specific criteria.
- Sample(s) D35708-1MS, D35708-1MSD, D35708-1SDL were used as the QC samples for the metals analysis.

## Metals By Method SW846 7471B

**Matrix** SO **Batch ID:** MP7755

- All samples were digested and analyzed within the recommended method holding time.
- All method blanks for this batch meet method specific criteria.
- Sample(s) D35710-1MS, D35710-1MSD were used as the QC samples for the metals analysis.

## Wet Chemistry By Method ASTM D1498-76M

**Matrix** SO **Batch ID:** GN15599

- Sample(s) D35794-1DUP were used as the QC samples for the Redox Potential Vs H2 analysis.

## Wet Chemistry By Method SM19 2540B M

**Matrix** SO **Batch ID:** GN15547

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

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

**Matrix** SO **Batch ID:** R13251

- The data for SW846 3060/7196A M meets quality control requirements.
- D35710-1 for Chromium, Trivalent: Calculated as: (Chromium) - (Chromium, Hexavalent)

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

<b>Matrix</b> SO	<b>Batch ID:</b> GP7549
------------------	-------------------------

- All samples were prepared and analyzed within the recommended method holding time.
- All method blanks for this batch meet method specific criteria.
- Sample(s) D35573-1DUP, D35573-1MS, D35573-1MSD were used as the QC samples for the Chromium, Hexavalent analysis.

### Wet Chemistry By Method SW846 9045D

<b>Matrix</b> SO	<b>Batch ID:</b> GN15564
------------------	--------------------------

- The following samples were run outside of holding time for method SW846 9045D: D35710-1

### Wet Chemistry By Method USDA HANDBOOK 60

<b>Matrix</b> SO	<b>Batch ID:</b> MP7743
------------------	-------------------------

- D35710-1A for Sodium Adsorption Ratio: Calculated as:  $(\text{Na meq/L}) / \sqrt{[(\text{Ca meq/L}) + (\text{Mg meq/L})/2]}$

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

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

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

Sample Results

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Report of Analysis

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## Report of Analysis

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<b>Client Sample ID:</b> CUT 1 SUBLINER	
<b>Lab Sample ID:</b> D35710-1	<b>Date Sampled:</b> 06/20/12
<b>Matrix:</b> SO - Soil	<b>Date Received:</b> 06/21/12
<b>Method:</b> SW846 8260B	<b>Percent Solids:</b> 92.1
<b>Project:</b> FRU 297-8B	

Run #	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	5V22147.D	1	06/25/12	BD	n/a	n/a	V5V1355
Run #2							

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

## Purgeable Aromatics

CAS No.	Compound	Result	RL	MDL	Units	Q
71-43-2	Benzene	ND	0.058	0.022	mg/kg	
108-88-3	Toluene	0.271	0.12	0.058	mg/kg	
100-41-4	Ethylbenzene	0.0912	0.12	0.022	mg/kg	J
1330-20-7	Xylene (total)	0.398	0.23	0.12	mg/kg	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
2037-26-5	Toluene-D8	93%		61-130%
460-00-4	4-Bromofluorobenzene	98%		53-131%
17060-07-0	1,2-Dichloroethane-D4	102%		62-130%

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

<b>Client Sample ID:</b> CUT 1 SUBLINER		
<b>Lab Sample ID:</b> D35710-1		<b>Date Sampled:</b> 06/20/12
<b>Matrix:</b> SO - Soil		<b>Date Received:</b> 06/21/12
<b>Method:</b> SW846 8270C BY SIM SW846 3546		<b>Percent Solids:</b> 92.1
<b>Project:</b> FRU 297-8B		

Run #	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	3G09824.D	1	06/22/12	SM	06/22/12	OP6113	E3G436
Run #2							

Run #	Initial Weight	Final Volume
Run #1	30.2 g	1.0 ml
Run #2		

## COGCC Table 910-1 PAH List

CAS No.	Compound	Result	RL	MDL	Units	Q
83-32-9	Acenaphthene	ND	0.0090	0.0047	mg/kg	
120-12-7	Anthracene	ND	0.0090	0.0047	mg/kg	
56-55-3	Benzo(a)anthracene	ND	0.0090	0.0047	mg/kg	
50-32-8	Benzo(a)pyrene	ND	0.0090	0.0047	mg/kg	
205-99-2	Benzo(b)fluoranthene	ND	0.0090	0.0047	mg/kg	
207-08-9	Benzo(k)fluoranthene	ND	0.0090	0.0047	mg/kg	
218-01-9	Chrysene	ND	0.0090	0.0047	mg/kg	
53-70-3	Dibenzo(a,h)anthracene	ND	0.0090	0.0047	mg/kg	
206-44-0	Fluoranthene	ND	0.0090	0.0047	mg/kg	
86-73-7	Fluorene	ND	0.0090	0.0047	mg/kg	
193-39-5	Indeno(1,2,3-cd)pyrene	ND	0.0090	0.0047	mg/kg	
91-20-3	Naphthalene	0.0549	0.013	0.011	mg/kg	
129-00-0	Pyrene	ND	0.0090	0.0047	mg/kg	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
4165-60-0	Nitrobenzene-d5	63%		10-145%
321-60-8	2-Fluorobiphenyl	78%		10-130%
1718-51-0	Terphenyl-d14	84%		22-130%

ND = Not detected MDL - Method Detection Limit

RL = Reporting Limit

E = Indicates value exceeds calibration range

J = Indicates an estimated value

B = Indicates analyte found in associated method blank

N = Indicates presumptive evidence of a compound

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## Report of Analysis

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<b>Client Sample ID:</b> CUT 1 SUBLINER	
<b>Lab Sample ID:</b> D35710-1	<b>Date Sampled:</b> 06/20/12
<b>Matrix:</b> SO - Soil	<b>Date Received:</b> 06/21/12
<b>Method:</b> SW846 8015B	<b>Percent Solids:</b> 92.1
<b>Project:</b> FRU 297-8B	

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	GB16420.D	1	06/21/12	SK	n/a	n/a	GGB910
Run #2							

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

CAS No.	Compound	Result	RL	MDL	Units	Q
	TPH-GRO (C6-C10)	ND	12	5.8	mg/kg	
CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits		
120-82-1	1,2,4-Trichlorobenzene	95%		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

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**Report of Analysis**

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<b>Client Sample ID:</b>	CUT 1 SUBLINER	<b>Date Sampled:</b>	06/20/12
<b>Lab Sample ID:</b>	D35710-1	<b>Date Received:</b>	06/21/12
<b>Matrix:</b>	SO - Soil	<b>Percent Solids:</b>	92.1
<b>Method:</b>	SW846-8015B SW846 3546		
<b>Project:</b>	FRU 297-8B		

	<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	FD14720.D	1	06/23/12	AW	06/22/12	OP6112	GFD764
Run #2							

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

<b>CAS No.</b>	<b>Compound</b>	<b>Result</b>	<b>RL</b>	<b>MDL</b>	<b>Units</b>	<b>Q</b>
	TPH-DRO (C10-C28)	83.9	14	9.4	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	83%		43-136%		

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

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

# Report of Analysis

<b>Client Sample ID:</b> CUT 1 SUBLINER	<b>Date Sampled:</b> 06/20/12
<b>Lab Sample ID:</b> D35710-1	<b>Date Received:</b> 06/21/12
<b>Matrix:</b> SO - Soil	<b>Percent Solids:</b> 92.1
<b>Project:</b> FRU 297-8B	

## Metals Analysis

Analyte	Result	RL	Units	DF	Prep	Analyzed By	Method	Prep Method
Arsenic	8.3	0.11	mg/kg	5	06/22/12	06/24/12 JM	SW846 6020A <sup>2</sup>	SW846 3050B <sup>5</sup>
Barium	533	1.1	mg/kg	1	06/22/12	06/23/12 JM	SW846 6010C <sup>1</sup>	SW846 3050B <sup>4</sup>
Cadmium	< 1.1	1.1	mg/kg	1	06/22/12	06/23/12 JM	SW846 6010C <sup>1</sup>	SW846 3050B <sup>4</sup>
Chromium	36.3	1.1	mg/kg	1	06/22/12	06/23/12 JM	SW846 6010C <sup>1</sup>	SW846 3050B <sup>4</sup>
Copper	13.1	1.1	mg/kg	1	06/22/12	06/23/12 JM	SW846 6010C <sup>1</sup>	SW846 3050B <sup>4</sup>
Lead	12.2	5.4	mg/kg	1	06/22/12	06/23/12 JM	SW846 6010C <sup>1</sup>	SW846 3050B <sup>4</sup>
Mercury	< 0.11	0.11	mg/kg	1	06/26/12	06/26/12 JM	SW846 7471B <sup>3</sup>	SW846 7471B <sup>6</sup>
Nickel	16.3	3.3	mg/kg	1	06/22/12	06/23/12 JM	SW846 6010C <sup>1</sup>	SW846 3050B <sup>4</sup>
Selenium	< 5.4	5.4	mg/kg	1	06/22/12	06/23/12 JM	SW846 6010C <sup>1</sup>	SW846 3050B <sup>4</sup>
Silver	< 3.3	3.3	mg/kg	1	06/22/12	06/23/12 JM	SW846 6010C <sup>1</sup>	SW846 3050B <sup>4</sup>
Zinc	46.5	3.3	mg/kg	1	06/22/12	06/23/12 JM	SW846 6010C <sup>1</sup>	SW846 3050B <sup>4</sup>

- (1) Instrument QC Batch: MA2541
- (2) Instrument QC Batch: MA2543
- (3) Instrument QC Batch: MA2547
- (4) Prep QC Batch: MP7737
- (5) Prep QC Batch: MP7738
- (6) Prep QC Batch: MP7755

RL = Reporting Limit

## Report of Analysis

<b>Client Sample ID:</b> CUT 1 SUBLINER	<b>Date Sampled:</b> 06/20/12
<b>Lab Sample ID:</b> D35710-1	<b>Date Received:</b> 06/21/12
<b>Matrix:</b> SO - Soil	<b>Percent Solids:</b> 92.1
<b>Project:</b> FRU 297-8B	

### General Chemistry

Analyte	Result	RL	Units	DF	Analyzed	By	Method
Chromium, Hexavalent	< 1.0	1.0	mg/kg	1	06/25/12	CJ	SW846 3060A/7196A
Chromium, Trivalent <sup>a</sup>	36.3	2.1	mg/kg	1	06/25/12	CJ	SW846 3060/7196A M
Redox Potential Vs H2	250		mv	1	06/26/12 12:50	JK	ASTM D1498-76M
Solids, Percent	92.1		%	1	06/22/12	SWT	SM19 2540B M
Specific Conductivity	8130	1.0	umhos/cm	1	06/22/12	CJ	DEPT.OF AG, BOOK N9
pH	10.49		su	1	06/22/12 13:15	JK	SW846 9045D

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

RL = Reporting Limit

## Report of Analysis

<b>Client Sample ID:</b> CUT 1 SUBLINER	<b>Date Sampled:</b> 06/20/12
<b>Lab Sample ID:</b> D35710-1A	<b>Date Received:</b> 06/21/12
<b>Matrix:</b> SO - Soil	<b>Percent Solids:</b> 92.1
<b>Project:</b> FRU 297-8B	

### SAR Metals Analysis

Analyte	Result	RL	Units	DF	Prep	Analyzed By	Method	Prep Method
Calcium	13.0	2.0	mg/l	1	06/22/12	06/23/12 JM	SW846 6010C <sup>1</sup>	EPA 200.7 <sup>2</sup>
Magnesium	7.51	1.0	mg/l	1	06/22/12	06/23/12 JM	SW846 6010C <sup>1</sup>	EPA 200.7 <sup>2</sup>
Sodium	1860	2.0	mg/l	1	06/22/12	06/23/12 JM	SW846 6010C <sup>1</sup>	EPA 200.7 <sup>2</sup>

(1) Instrument QC Batch: MA2541

(2) Prep QC Batch: MP7743

---

RL = Reporting Limit

## Report of Analysis

<b>Client Sample ID:</b> CUT 1 SUBLINER	<b>Date Sampled:</b> 06/20/12
<b>Lab Sample ID:</b> D35710-1A	<b>Date Received:</b> 06/21/12
<b>Matrix:</b> SO - Soil	<b>Percent Solids:</b> 92.1
<b>Project:</b> FRU 297-8B	

### General Chemistry

Analyte	Result	RL	Units	DF	Analyzed	By	Method
Sodium Adsorption Ratio <sup>a</sup>	102		ratio	1	06/23/12 16:19	JM	USDA HANDBOOK 60

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

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RL = Reporting Limit

Misc. Forms

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Custody Documents and Other Forms

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Includes the following where applicable:

- Chain of Custody



CHAIN OF CUSTODY

4035 Youngfield Street, Wheat Ridge, CO 80033
TEL: 303-425-6021 FAX: 303-425-6834
www.accutest.com

FED-EX Tracking #
Bottle Order Control #
Accutest Quote #
Accutest Job # D35710

Client / Reporting Information
Project Information
Requested Analysis (see TEST CODE sheet)
Matrix Codes
Company Name: KRW Consulting
Project Name: XTO FRU 297-8B
Street Address: 8000 West 14th Street, Suite 200
City: Lakewood, CO 80214
Project Contact: Dwayne Knudson
Project # 1106-06
Client Purchase Order #
City: Rifle, CO 81650
Sampler(s) Name(s): DAVID SANDERS
Project Manager: Joe Hess
Attention: Jessica Dooling

Table with columns: Turnaround Time (Business days), Data Deliverable Information, Comments / Special Instructions. Includes checkboxes for Std. 10 Business Days, Std. 6 Business Days, 6 Day RUSH, 3 Day Emergency, 2 Day Emergency, 1 Day Emergency. Also includes checkboxes for Commercial 'A' (Level 1), Commercial 'B' (Level 2), COMMEN, COMMEN+, State Forms Required, Send Forms to State, Report by Fax, Report by PDF ONLY, EDD Format.

Sample Custody must be documented below each time samples change possession, including courier delivery.
Relinquished by: 1. David Sanders Date Time: 6/20/12 3:50
Received By: 1. [Signature] Date Time: 6/20/12
Relinquished by: 2. [Signature] Date Time: [Signature]
Received By: 2. [Signature] Date Time: [Signature]
Relinquished by: 3. [Signature] Date Time: [Signature]
Received By: 3. [Signature] Date Time: [Signature]
Relinquished by: 4. [Signature] Date Time: [Signature]
Received By: 4. [Signature] Date Time: [Signature]
Relinquished by: 5. [Signature] Date Time: [Signature]
Received By: 5. [Signature] Date Time: [Signature]
Relinquished by: 6. [Signature] Date Time: [Signature]
Received By: 6. [Signature] Date Time: [Signature]

4.1
4

D35710: Chain of Custody

Page 1 of 2

Accutest Job Number: D35710

Client: KRW

Immediate Client Services Action Required: No

Date / Time Received: 6/21/2012 10:20:00 AM

No. Coolers: 1

Client Service Action Required at Login: No

Project: XTO

Airbill #'s: CO

<u>Cooler Security</u>	<u>Y or N</u>		<u>Y or N</u>	
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"/>

<u>Cooler Temperature</u>	<u>Y or N</u>	
1. Temp criteria achieved:	<input checked="" type="checkbox"/>	<input type="checkbox"/>
2. Cooler temp verification:	Infrared gun	
3. Cooler media:	Ice (bag)	

<u>Quality Control Preservation</u>	<u>Y or N</u>		<u>N/A</u>
1. Trip Blank present / cooler:	<input type="checkbox"/>	<input type="checkbox"/>	
2. Trip Blank listed on COC:	<input type="checkbox"/>	<input type="checkbox"/>	
3. Samples preserved properly:	<input checked="" type="checkbox"/>	<input type="checkbox"/>	
4. VOCs headspace free:	<input type="checkbox"/>	<input type="checkbox"/>	<input checked="" type="checkbox"/>

<u>Sample Integrity - Documentation</u>	<u>Y or N</u>	
1. Sample labels present on bottles:	<input checked="" type="checkbox"/>	<input type="checkbox"/>
2. Container labeling complete:	<input checked="" type="checkbox"/>	<input type="checkbox"/>
3. Sample container label / COC agree:	<input checked="" type="checkbox"/>	<input type="checkbox"/>

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

<u>Sample Integrity - Instructions</u>	<u>Y or N</u>		<u>N/A</u>
1. Analysis requested is clear:	<input checked="" type="checkbox"/>	<input type="checkbox"/>	
2. Bottles received for unspecified tests	<input type="checkbox"/>	<input checked="" type="checkbox"/>	
3. Sufficient volume rec'd for analysis:	<input checked="" type="checkbox"/>	<input type="checkbox"/>	
4. Compositing instructions clear:	<input type="checkbox"/>	<input type="checkbox"/>	<input checked="" type="checkbox"/>
5. Filtering instructions clear:	<input type="checkbox"/>	<input type="checkbox"/>	<input checked="" type="checkbox"/>

Comments

4.1  
4

## GC/MS Volatiles

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5

## 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:** D35710  
**Account:** XTOKRWR XTO Energy  
**Project:** FRU 297-8B

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
V5V1355-MB	5V22134.D	1	06/25/12	BD	n/a	n/a	V5V1355

The QC reported here applies to the following samples:

Method: SW846 8260B

D35710-1

CAS No.	Compound	Result	RL	MDL	Units	Q
71-43-2	Benzene	ND	50	19	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	88%	61-130%
460-00-4	4-Bromofluorobenzene	88%	53-131%
17060-07-0	1,2-Dichloroethane-D4	110%	62-130%

# Blank Spike Summary

**Job Number:** D35710  
**Account:** XTOKRWR XTO Energy  
**Project:** FRU 297-8B

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
V5V1355-BS	5V22135.D	1	06/25/12	BD	n/a	n/a	V5V1355

The QC reported here applies to the following samples:

Method: SW846 8260B

D35710-1

CAS No.	Compound	Spike ug/kg	BSP ug/kg	BSP %	Limits
71-43-2	Benzene	50	49.0	98	70-130
100-41-4	Ethylbenzene	50	46.0	92	70-130
108-88-3	Toluene	50	43.2	86	70-130
1330-20-7	Xylene (total)	150	141	94	70-130

CAS No.	Surrogate Recoveries	BSP	Limits
2037-26-5	Toluene-D8	83%	61-130%
460-00-4	4-Bromofluorobenzene	93%	53-131%
17060-07-0	1,2-Dichloroethane-D4	98%	62-130%

\* = Outside of Control Limits.

# Matrix Spike/Matrix Spike Duplicate Summary

**Job Number:** D35710  
**Account:** XTOKRWR XTO Energy  
**Project:** FRU 297-8B

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
D35708-1MS	5V22137.D	1	06/25/12	BD	n/a	n/a	V5V1355
D35708-1MSD	5V22138.D	1	06/25/12	BD	n/a	n/a	V5V1355
D35708-1	5V22136.D	1	06/25/12	BD	n/a	n/a	V5V1355

The QC reported here applies to the following samples:

Method: SW846 8260B

D35710-1

CAS No.	Compound	D35708-1 ug/kg	Q	Spike ug/kg	MS ug/kg	MS %	MSD ug/kg	MSD %	RPD	Limits Rec/RPD
71-43-2	Benzene	22.0	J	2750	2880	104	2870	104	0	70-134/30
100-41-4	Ethylbenzene	149		2750	2770	95	2740	94	1	70-137/30
108-88-3	Toluene	445		2750	2740	83	2740	83	0	70-130/30
1330-20-7	Xylene (total)	614		8250	8670	98	8540	96	2	61-131/30

CAS No.	Surrogate Recoveries	MS	MSD	D35708-1	Limits
2037-26-5	Toluene-D8	88%	85%	95%	61-130%
460-00-4	4-Bromofluorobenzene	108%	103%	103%	53-131%
17060-07-0	1,2-Dichloroethane-D4	107%	96%	114%	62-130%

\* = Outside of Control Limits.

5.3.1  
 5

GC/MS Volatiles

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Raw Data

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## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\V5062512.S\  
 Data File : 5V22147.D  
 Acq On : 25 Jun 2012 7:38 pm  
 Operator : BRETD  
 Sample : D35710-1, X50  
 Misc : MS4172,V5V1355,5.044,,100,5,1  
 ALS Vial : 16 Sample Multiplier: 1

Quant Time: Jun 26 09:03:12 2012  
 Quant Method : C:\msdchem\1\METHODS\V5AP1304TVH1304.M  
 Quant Title : 8260  
 QLast Update : Thu May 24 07:55:17 2012  
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
2) Pentafluorobenzene	11.647	168	178632	50.00	ug/l	0.00
35) 1,4-Difluorobenzene	12.446	114	284115	50.00	ug/l	0.00
53) Chlorobenzene-d5	15.095	117	367244	50.00	ug/l	0.00
74) 1,4-Dichlorobenzene-d4	17.070	152	259467	50.00	ug/l	0.00

## System Monitoring Compounds

33) 1,2-Dichloroethane-d4	12.035	102	27996	51.17	ug/l	0.00
Spiked Amount	50.000	Range	70 - 130	Recovery	=	102.34%
61) Toluene-d8	13.850	98	576899	46.48	ug/l	0.00
Spiked Amount	50.000	Range	70 - 130	Recovery	=	92.96%
69) 4-Bromofluorobenzene	16.043	95	249671	49.11	ug/l	0.00
Spiked Amount	50.000	Range	70 - 130	Recovery	=	98.22%

## Target Compounds

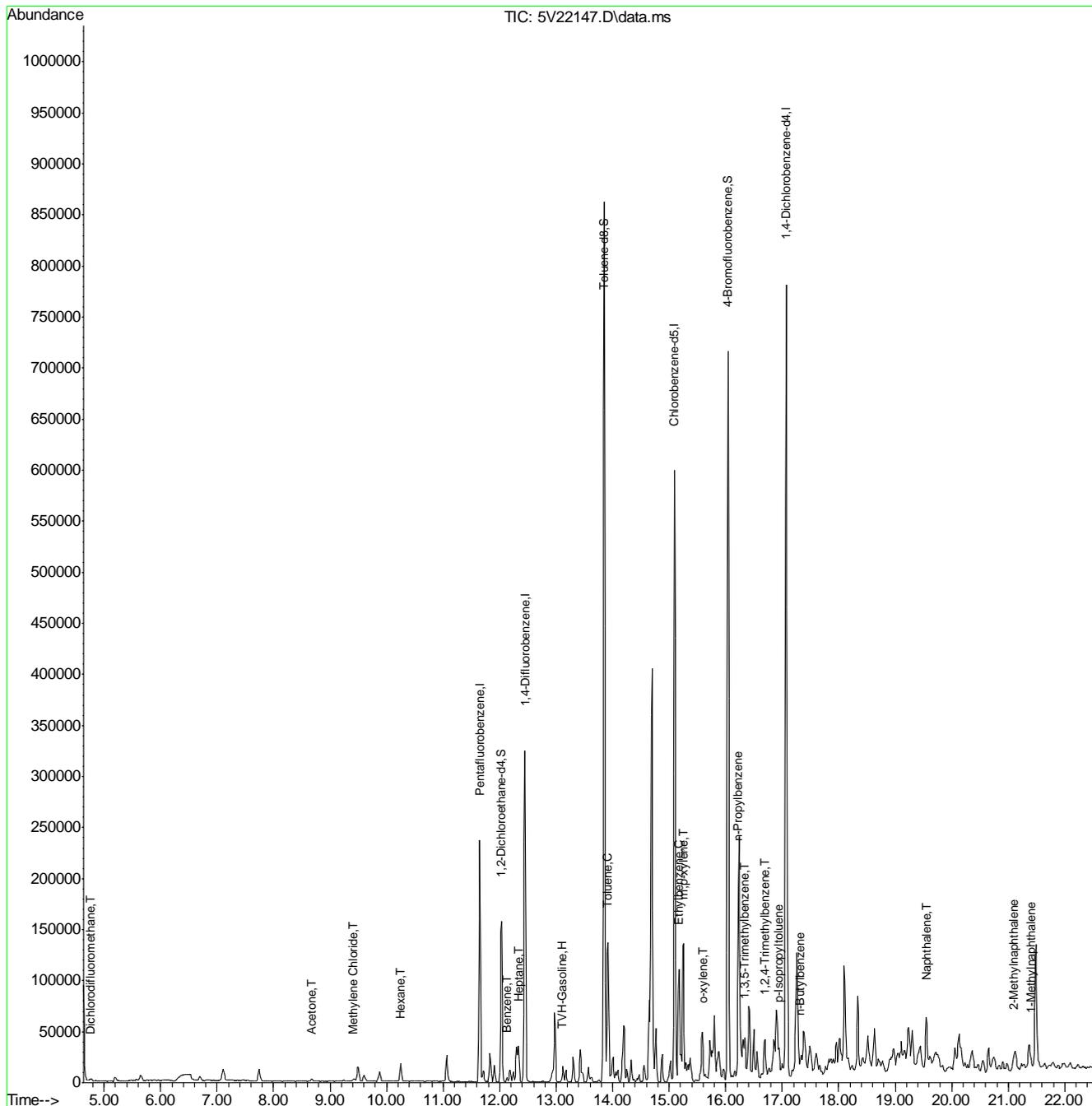
						Qvalue
1) TVH-Gasoline	13.102	TIC	4065616m	204.18	ug/l	
3) Dichlorodifluoromethane	4.763	85	2065	1.13	ug/l	74
15) Acetone	8.667	58	790	0.89	ug/l #	62
17) Methylene Chloride	9.409	84	1477	0.49	ug/l	85
41) Hexane	10.243	57	9369	2.38	ug/l	100
43) Heptane	12.332	43	13725	3.26	ug/l	90
50) Benzene	12.127	78	2548	0.22	ug/l	100
62) Toluene	13.908	92	44530	4.66	ug/l	99
66) Ethylbenzene	15.175	91	27772	1.57	ug/l	99
72) m,p-xylene	15.255	106	43882	6.20	ug/l	97
73) o-xylene	15.597	106	4378	0.64	ug/l	88
77) n-Propylbenzene	16.225	91	10268	0.44	ug/l	96
80) 1,3,5-Trimethylbenzene	16.328	105	2207m	0.13	ug/l	
82) 1,2,4-Trimethylbenzene	16.693	105	14122	0.84	ug/l	89
86) p-Isopropyltoluene	16.944	119	13177	0.70	ug/l	96
88) n-Butylbenzene	17.321	91	7925	0.46	ug/l #	75
91) Naphthalene	19.559	128	10872	1.40	ug/l	100
94) 2-Methylnaphthalene	21.100	142	10686	3.31	ug/l #	90
95) 1-Methylnaphthalene	21.397	142	4842	2.09	ug/l #	80

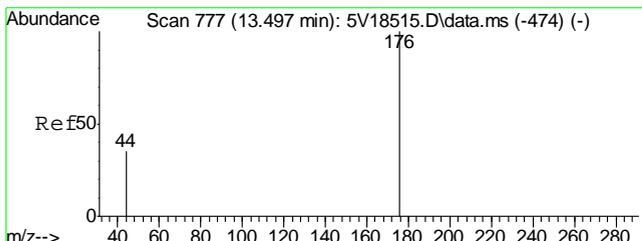
(#) = qualifier out of range (m) = manual integration (+) = signals summed

Quantitation Report (QT Reviewed)

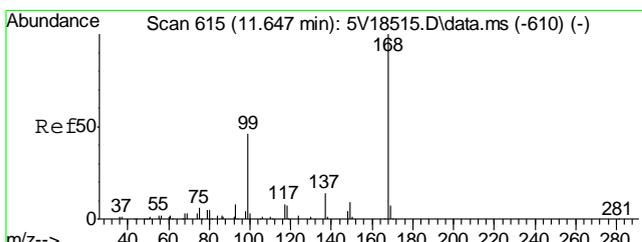
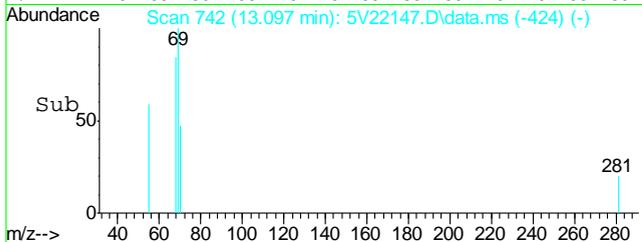
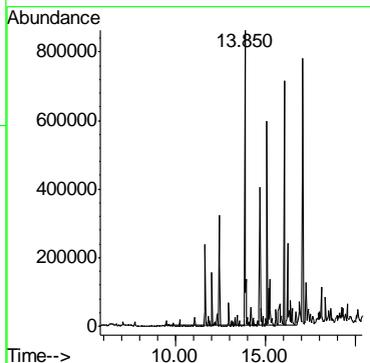
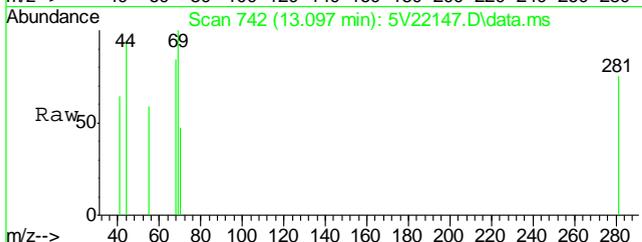
Data Path : C:\msdchem\1\DATA\V5062512.S\  
 Data File : 5V22147.D  
 Acq On : 25 Jun 2012 7:38 pm  
 Operator : BRETD  
 Sample : D35710-1, X50  
 Misc : MS4172,V5V1355,5.044,,100,5,1  
 ALS Vial : 16 Sample Multiplier: 1

Quant Time: Jun 26 09:03:12 2012  
 Quant Method : C:\msdchem\1\METHODS\V5AP1304TVH1304.M  
 Quant Title : 8260  
 QLast Update : Thu May 24 07:55:17 2012  
 Response via : Initial Calibration

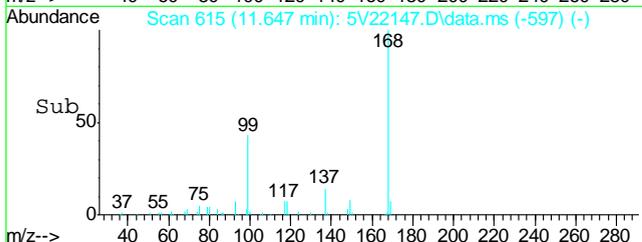
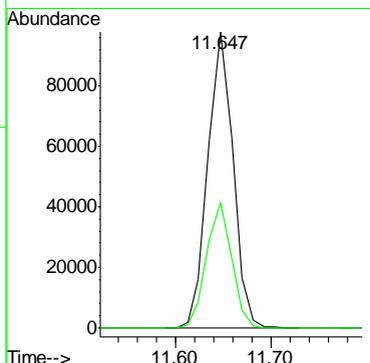
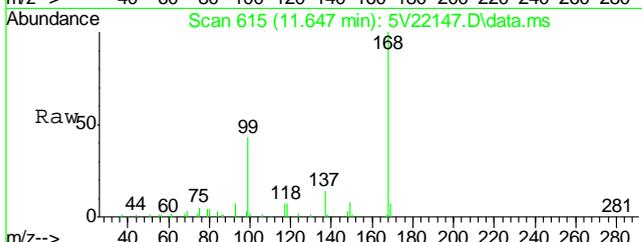


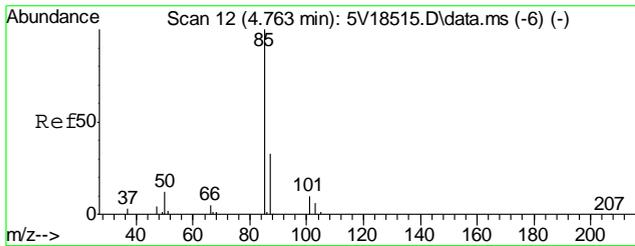


#1  
 TVH-Gasoline  
 Concen: 204.18 ug/l m  
 RT: 13.102 min Scan# 742  
 Delta R.T. 0.000 min  
 Lab File: 5V22147.D  
 Acq: 25 Jun 2012 7:38 pm  
 Tgt Ion:TIC Resp: 4065616



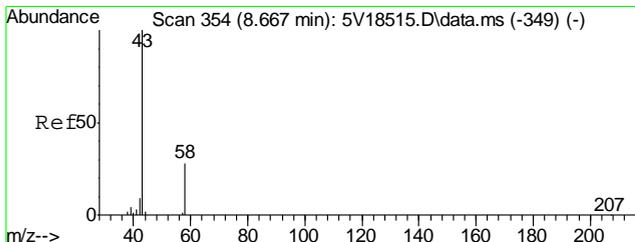
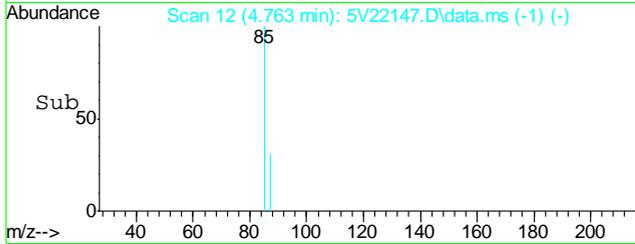
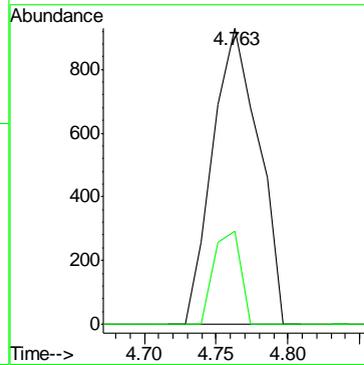
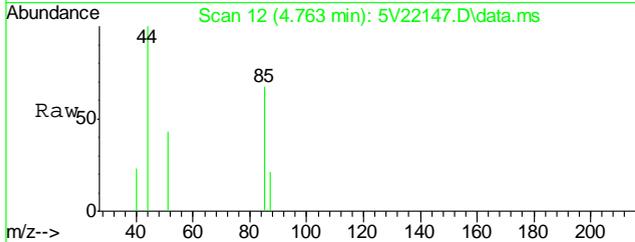
#2  
 Pentafluorobenzene  
 Concen: 50.00 ug/l  
 RT: 11.647 min Scan# 615  
 Delta R.T. -0.000 min  
 Lab File: 5V22147.D  
 Acq: 25 Jun 2012 7:38 pm  
 Tgt Ion:168 Resp: 178632  
 Ion Ratio Lower Upper  
 168 100  
 99 42.3 37.4 56.2





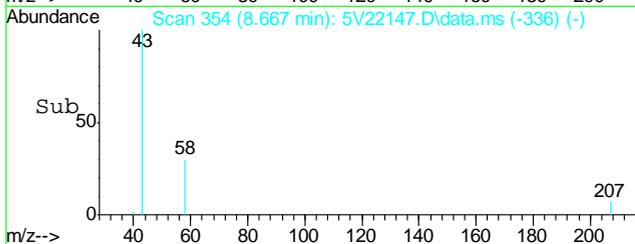
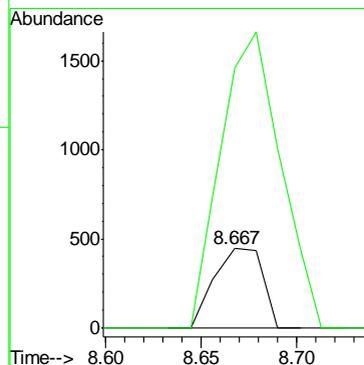
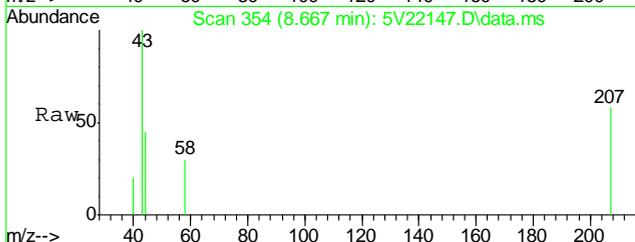
#3  
 Dichlorodifluoromethane  
 Concen: 1.13 ug/l  
 RT: 4.763 min Scan# 12  
 Delta R.T. 0.001 min  
 Lab File: 5V22147.D  
 Acq: 25 Jun 2012 7:38 pm

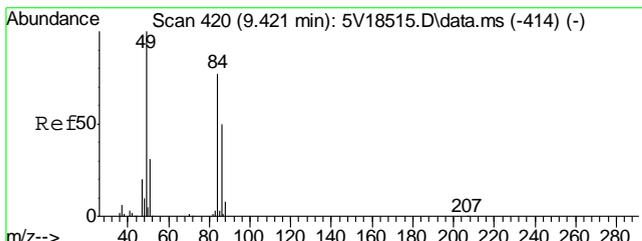
Tgt Ion	Resp	Lower	Upper
85	2065		
85	100		
87	18.2	12.9	52.9



#15  
 Acetone  
 Concen: 0.89 ug/l  
 RT: 8.667 min Scan# 354  
 Delta R.T. 0.000 min  
 Lab File: 5V22147.D  
 Acq: 25 Jun 2012 7:38 pm

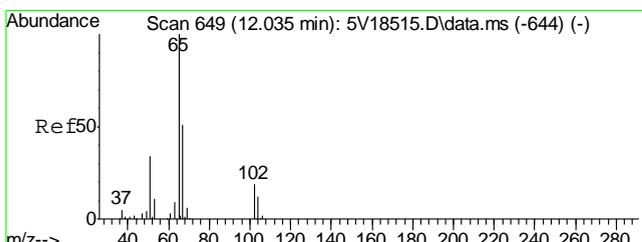
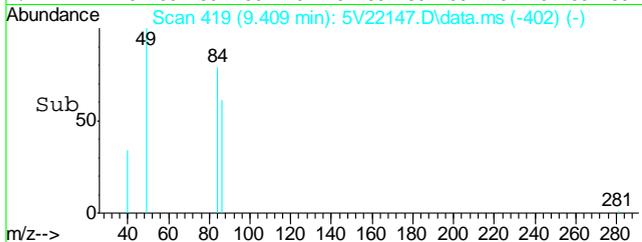
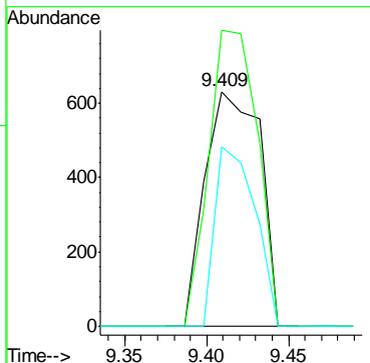
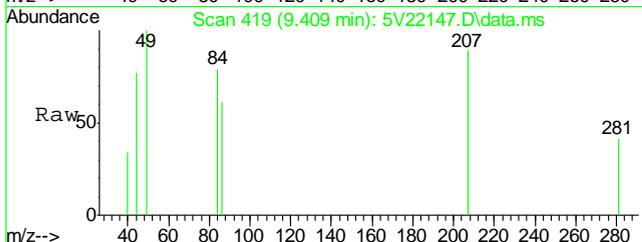
Tgt Ion	Resp	Lower	Upper
58	790		
58	100		
43	460.6	353.6	393.6#





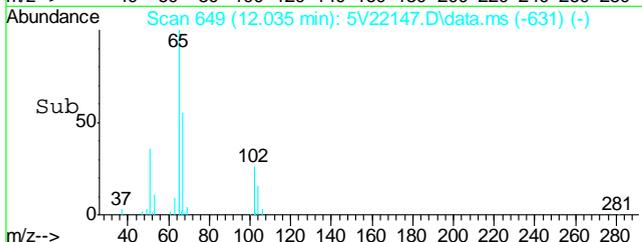
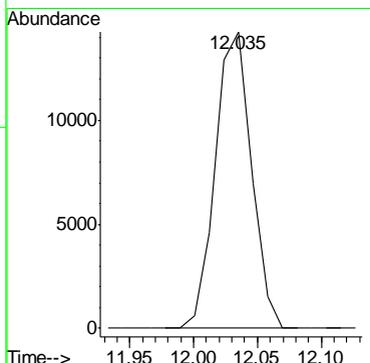
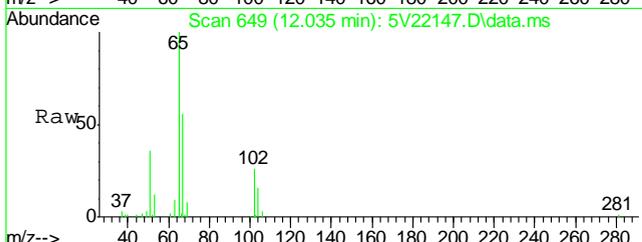
#17  
Methylene Chloride  
Concen: 0.49 ug/l  
RT: 9.409 min Scan# 419  
Delta R.T. -0.011 min  
Lab File: 5V22147.D  
Acq: 25 Jun 2012 7:38 pm

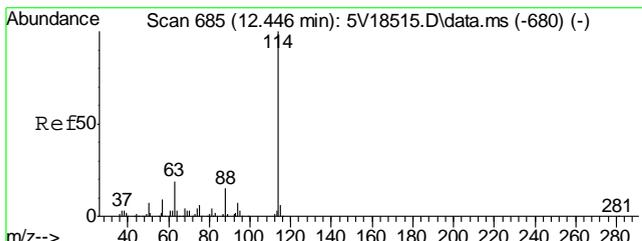
Tgt Ion	Resp	Lower	Upper
84	1477		
84	100		
49	110.8	110.4	150.4
86	55.5	44.0	84.0



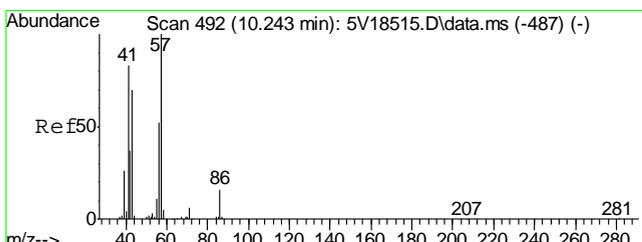
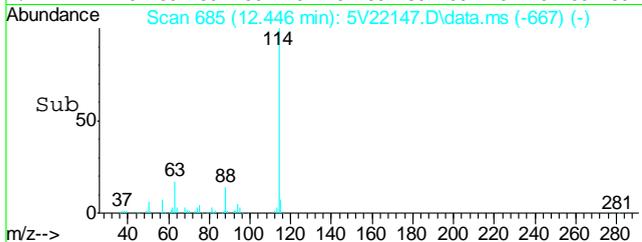
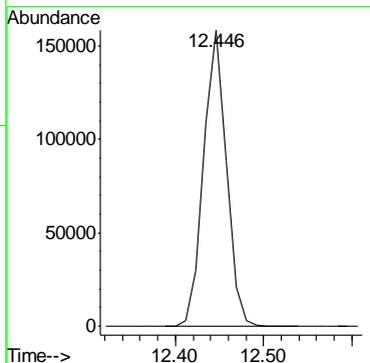
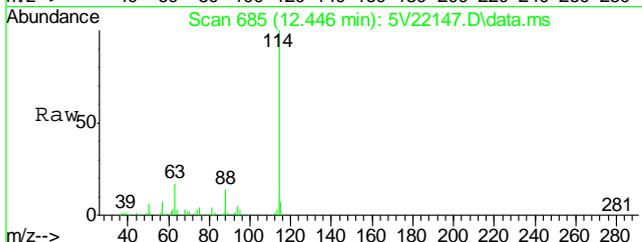
#33  
1,2-Dichloroethane-d4  
Concen: 51.17 ug/l  
RT: 12.035 min Scan# 649  
Delta R.T. 0.000 min  
Lab File: 5V22147.D  
Acq: 25 Jun 2012 7:38 pm

Tgt Ion	Resp
102	27996

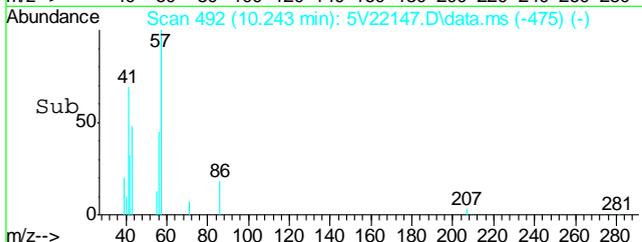
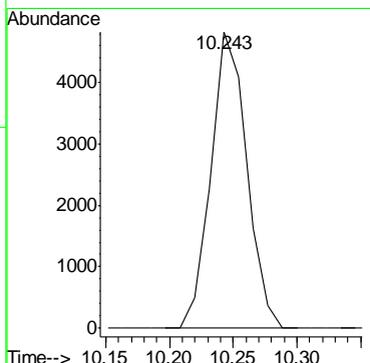
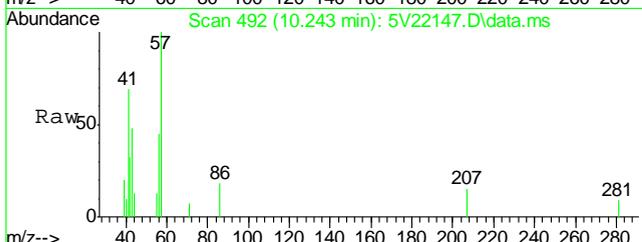


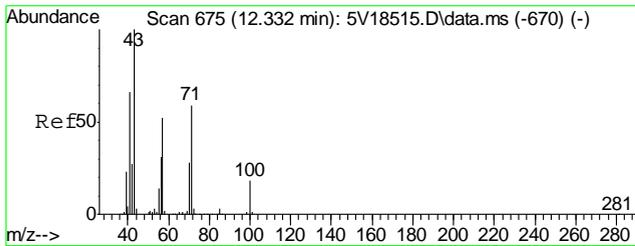


#35  
 1,4-Difluorobenzene  
 Concen: 50.00 ug/l  
 RT: 12.446 min Scan# 685  
 Delta R.T. -0.000 min  
 Lab File: 5V22147.D  
 Acq: 25 Jun 2012 7:38 pm  
 Tgt Ion:114 Resp: 284115



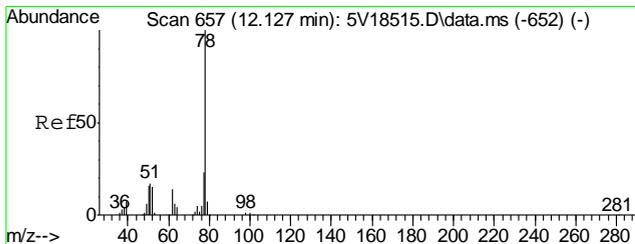
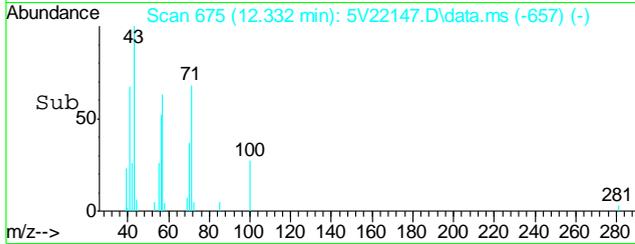
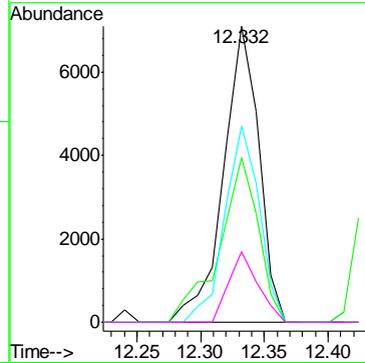
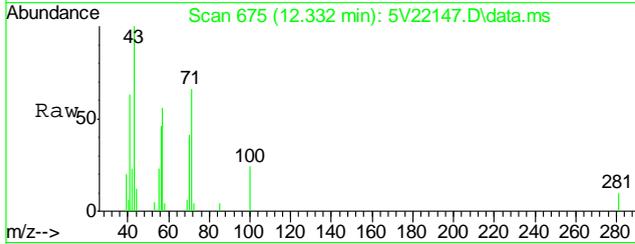
#41  
 Hexane  
 Concen: 2.38 ug/l  
 RT: 10.243 min Scan# 492  
 Delta R.T. -0.012 min  
 Lab File: 5V22147.D  
 Acq: 25 Jun 2012 7:38 pm  
 Tgt Ion: 57 Resp: 9369





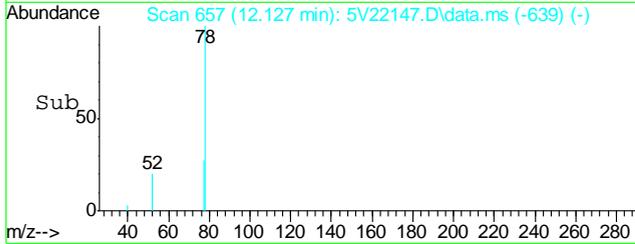
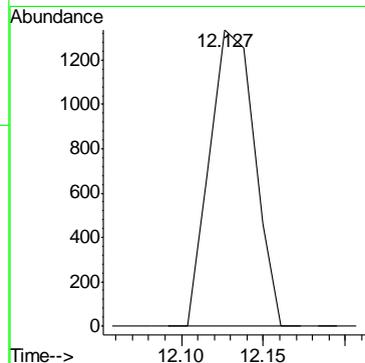
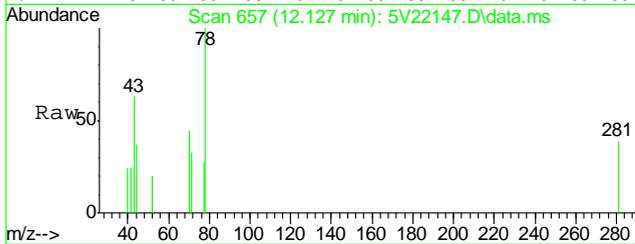
#43  
 Heptane  
 Concen: 3.26 ug/l  
 RT: 12.332 min Scan# 675  
 Delta R.T. -0.000 min  
 Lab File: 5V22147.D  
 Acq: 25 Jun 2012 7:38 pm

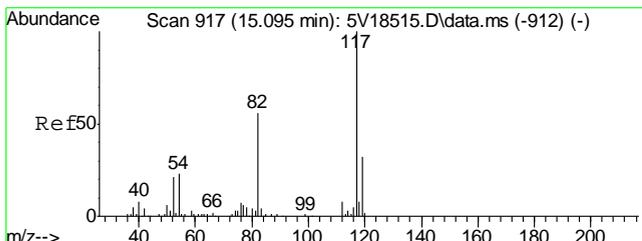
Tgt Ion	Resp	Lower	Upper
43	13725		
57	60.7	30.6	70.6
71	64.5	38.9	78.9
100	19.7	0.0	37.4



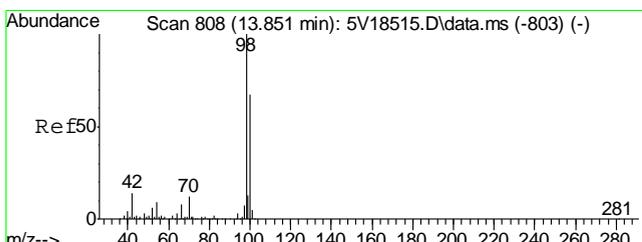
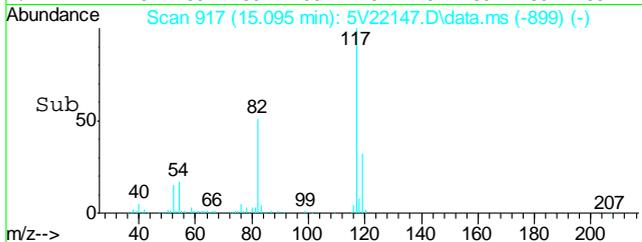
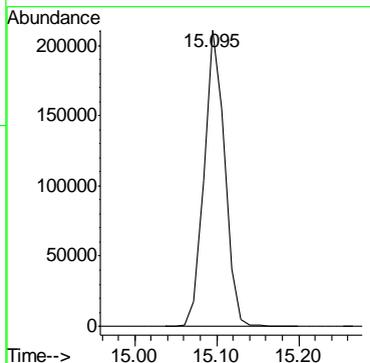
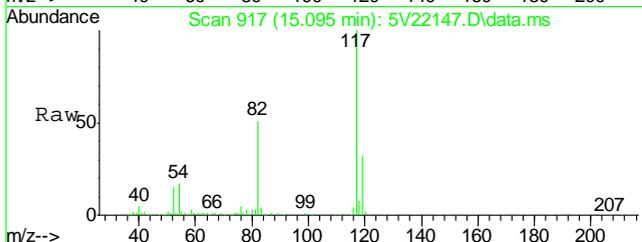
#50  
 Benzene  
 Concen: 0.22 ug/l  
 RT: 12.127 min Scan# 657  
 Delta R.T. -0.000 min  
 Lab File: 5V22147.D  
 Acq: 25 Jun 2012 7:38 pm

Tgt Ion: 78 Resp: 2548

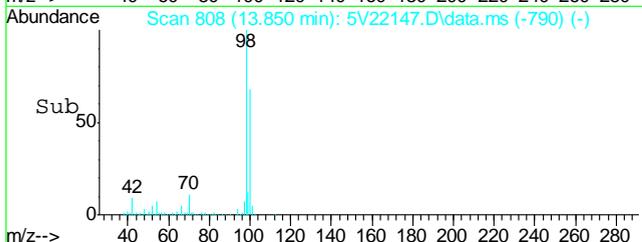
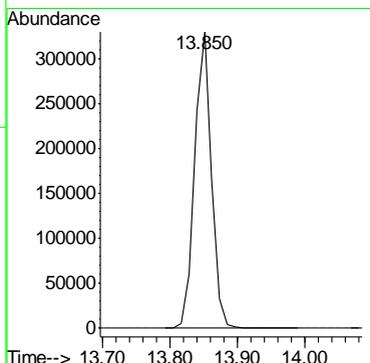
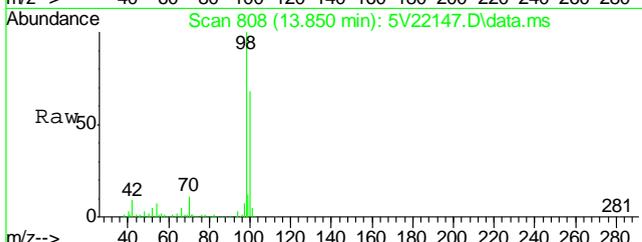


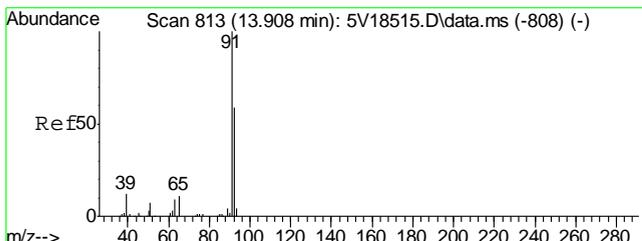


#53  
 Chlorobenzene-d5  
 Concen: 50.00 ug/l  
 RT: 15.095 min Scan# 917  
 Delta R.T. -0.000 min  
 Lab File: 5V22147.D  
 Acq: 25 Jun 2012 7:38 pm  
 Tgt Ion:117 Resp: 367244



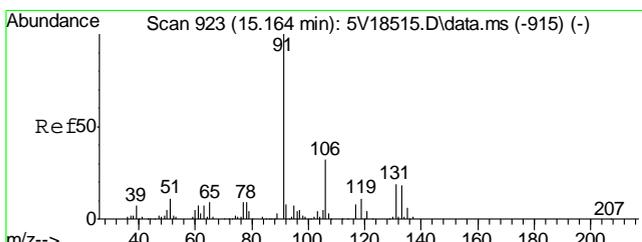
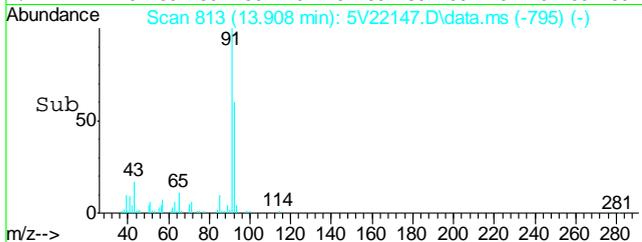
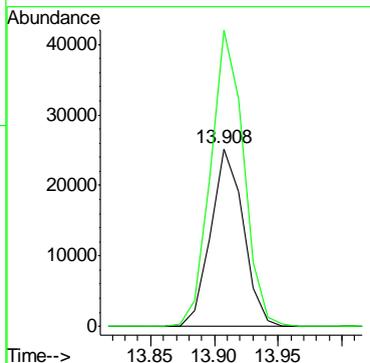
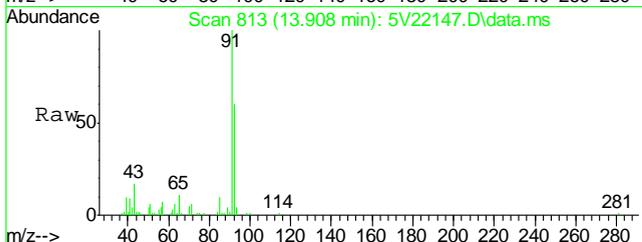
#61  
 Toluene-d8  
 Concen: 46.48 ug/l  
 RT: 13.850 min Scan# 808  
 Delta R.T. -0.000 min  
 Lab File: 5V22147.D  
 Acq: 25 Jun 2012 7:38 pm  
 Tgt Ion: 98 Resp: 576899





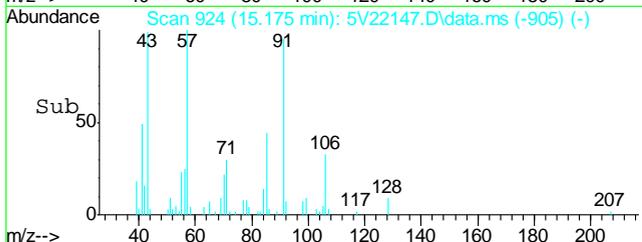
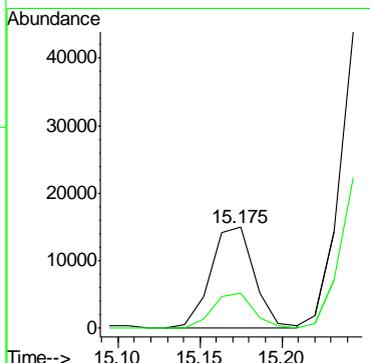
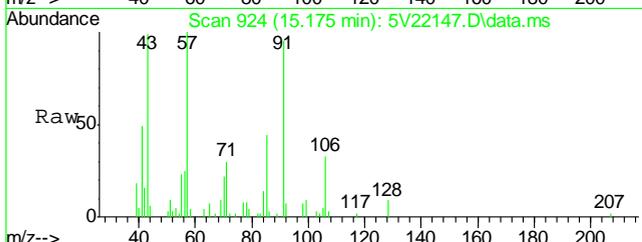
#62  
Toluene  
Concen: 4.66 ug/l  
RT: 13.908 min Scan# 813  
Delta R.T. -0.000 min  
Lab File: 5V22147.D  
Acq: 25 Jun 2012 7:38 pm

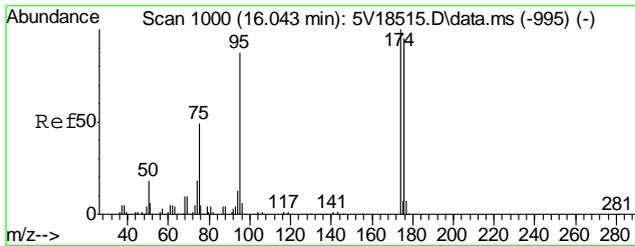
Tgt Ion	Resp	Lower	Upper
92	44530	100	
91	168.6	149.8	189.8



#66  
Ethylbenzene  
Concen: 1.57 ug/l  
RT: 15.175 min Scan# 924  
Delta R.T. 0.011 min  
Lab File: 5V22147.D  
Acq: 25 Jun 2012 7:38 pm

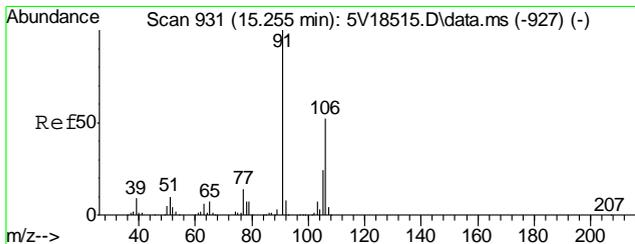
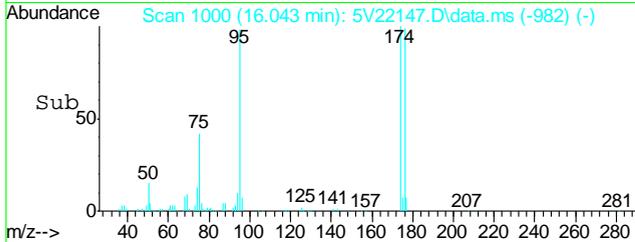
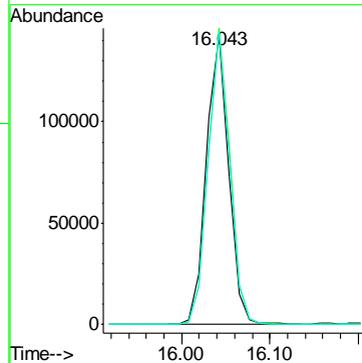
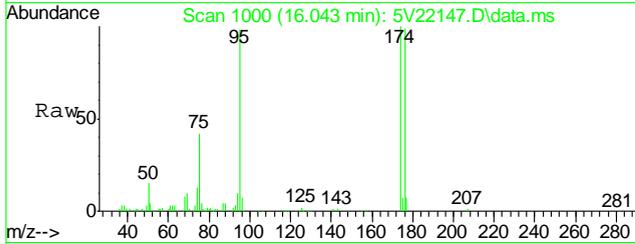
Tgt Ion	Resp	Lower	Upper
91	27772	100	
106	32.3	11.7	51.7





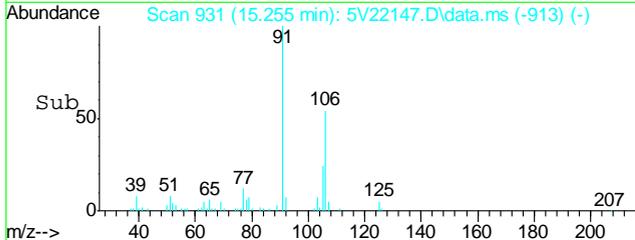
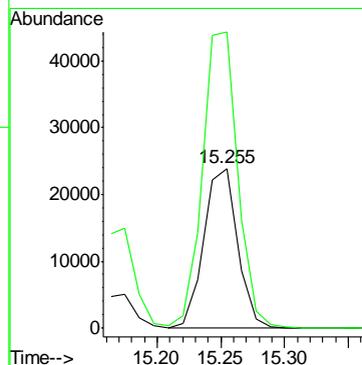
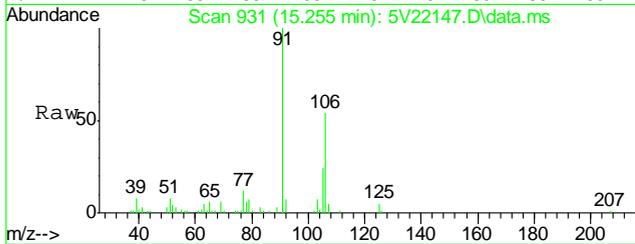
#69  
4-Bromofluorobenzene  
Concen: 49.11 ug/l  
RT: 16.043 min Scan# 1000  
Delta R.T. -0.000 min  
Lab File: 5V22147.D  
Acq: 25 Jun 2012 7:38 pm

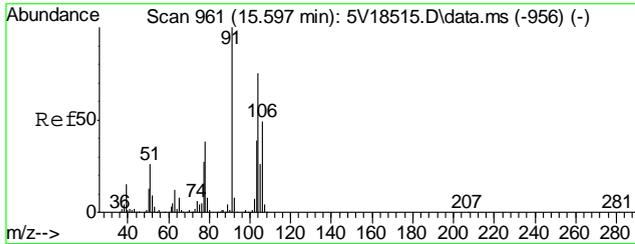
Tgt Ion	Resp	Lower	Upper
95	249671		
174	100.1	77.1	117.1
176	98.2	73.4	113.4



#72  
m,p-xylene  
Concen: 6.20 ug/l  
RT: 15.255 min Scan# 931  
Delta R.T. -0.000 min  
Lab File: 5V22147.D  
Acq: 25 Jun 2012 7:38 pm

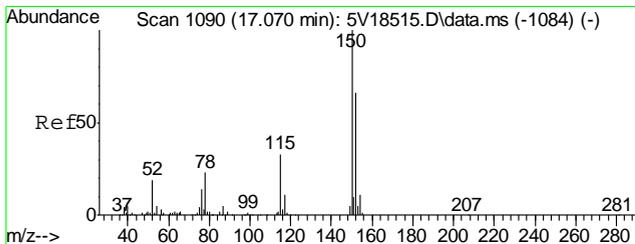
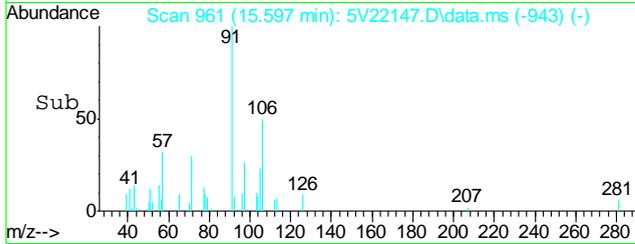
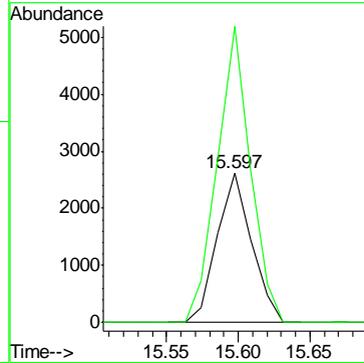
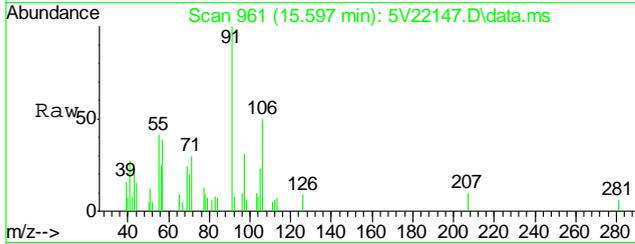
Tgt Ion	Resp	Lower	Upper
106	43882		
91	193.2	177.1	217.1





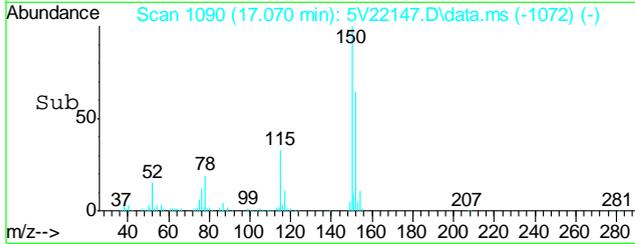
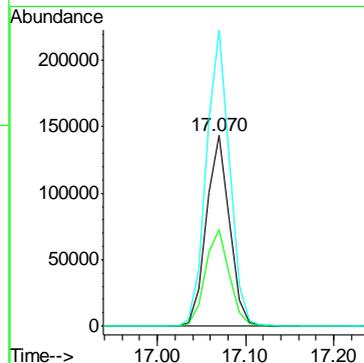
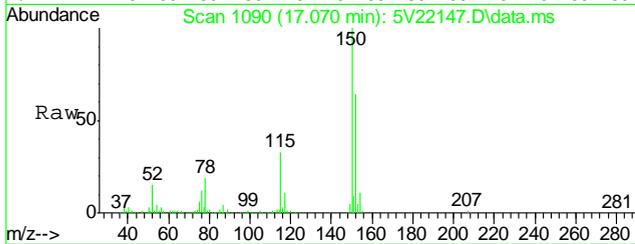
#73  
o-xylene  
Concen: 0.64 ug/l  
RT: 15.597 min Scan# 961  
Delta R.T. -0.000 min  
Lab File: 5V22147.D  
Acq: 25 Jun 2012 7:38 pm

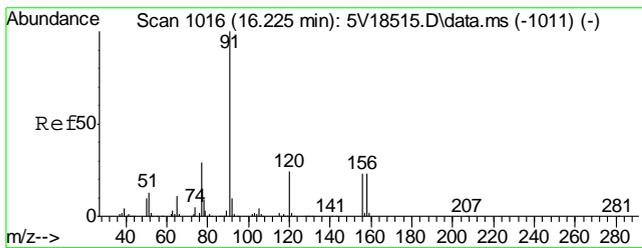
Tgt Ion	Resp	Lower	Upper
106	4378	100	
91	190.3	166.6	249.8



#74  
1,4-Dichlorobenzene-d4  
Concen: 50.00 ug/l  
RT: 17.070 min Scan# 1090  
Delta R.T. -0.000 min  
Lab File: 5V22147.D  
Acq: 25 Jun 2012 7:38 pm

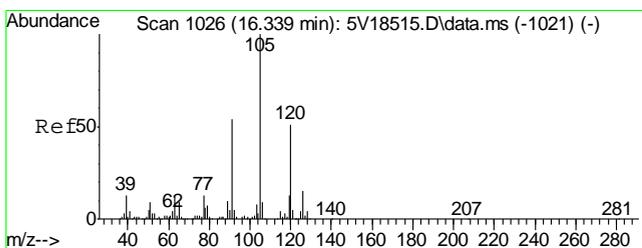
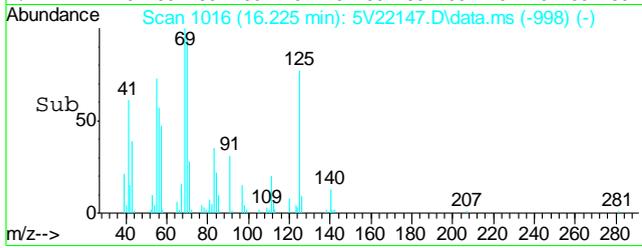
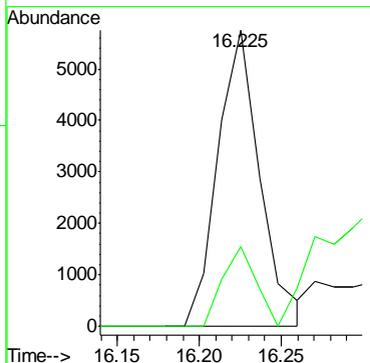
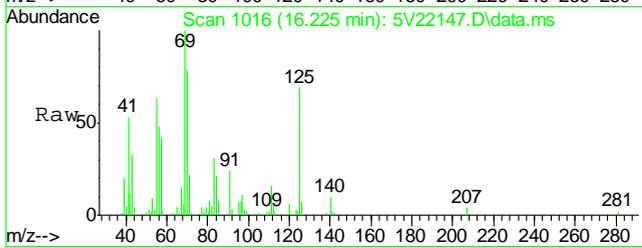
Tgt Ion	Resp	Lower	Upper
152	259467	100	
115	52.2	41.4	62.0
150	155.8	153.9	230.9





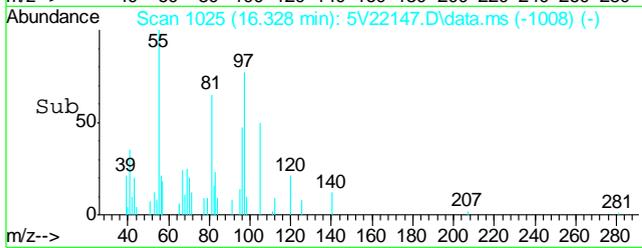
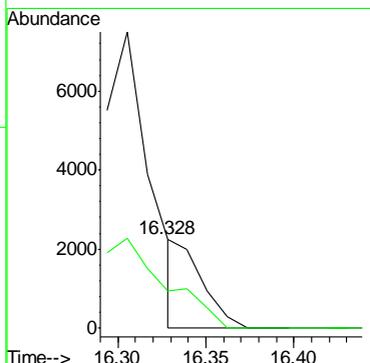
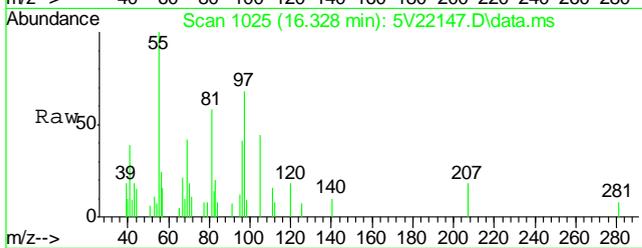
#77  
 n-Propylbenzene  
 Concen: 0.44 ug/l  
 RT: 16.225 min Scan# 1016  
 Delta R.T. -0.000 min  
 Lab File: 5V22147.D  
 Acq: 25 Jun 2012 7:38 pm

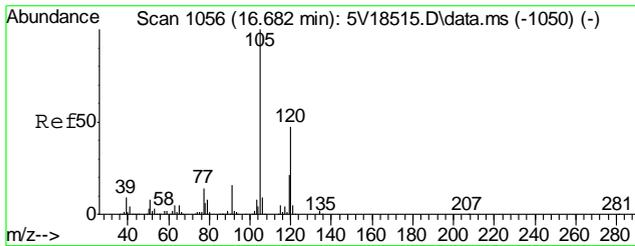
Tgt Ion	Resp	Lower	Upper
91	10268	100	
120	21.3	18.6	27.8



#80  
 1,3,5-Trimethylbenzene  
 Concen: 0.13 ug/l m  
 RT: 16.328 min Scan# 1025  
 Delta R.T. -0.011 min  
 Lab File: 5V22147.D  
 Acq: 25 Jun 2012 7:38 pm

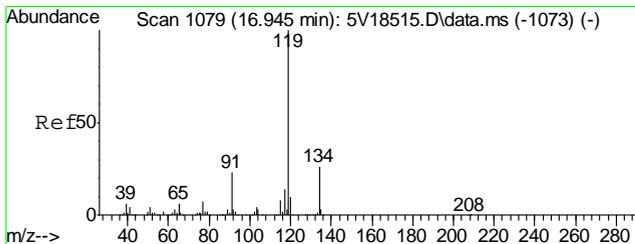
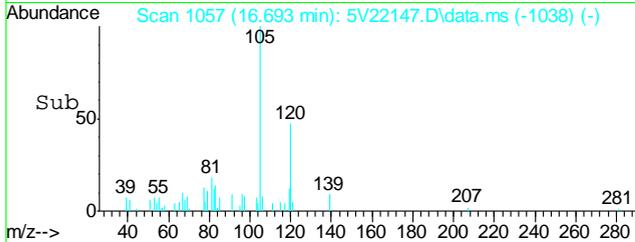
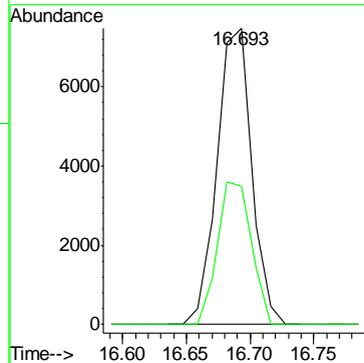
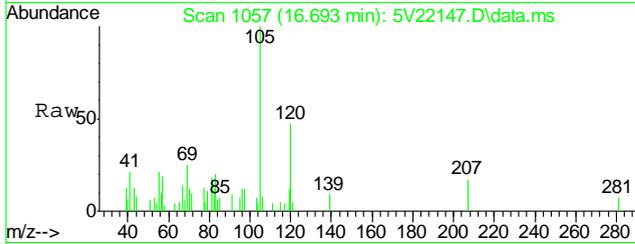
Tgt Ion	Resp	Lower	Upper
105	2207	100	
120	379.3	40.1	60.1#





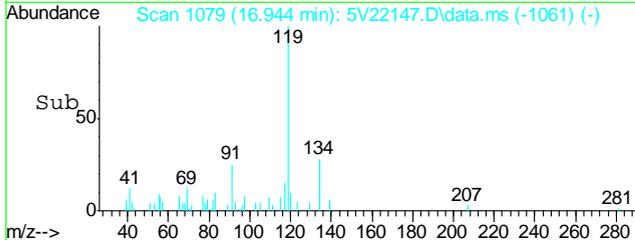
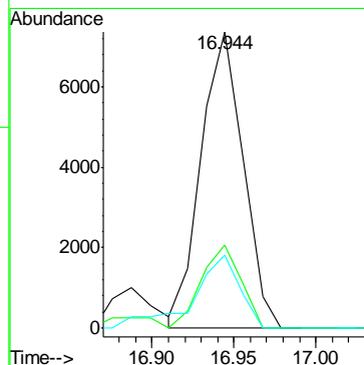
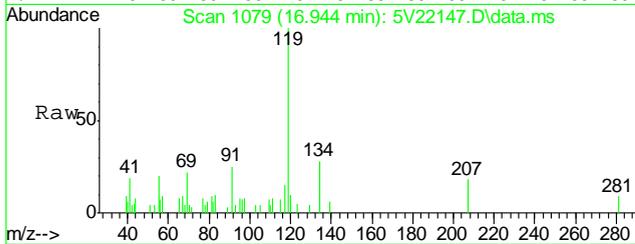
#82  
 1,2,4-Trimethylbenzene  
 Concen: 0.84 ug/l  
 RT: 16.693 min Scan# 1057  
 Delta R.T. 0.011 min  
 Lab File: 5V22147.D  
 Acq: 25 Jun 2012 7:38 pm

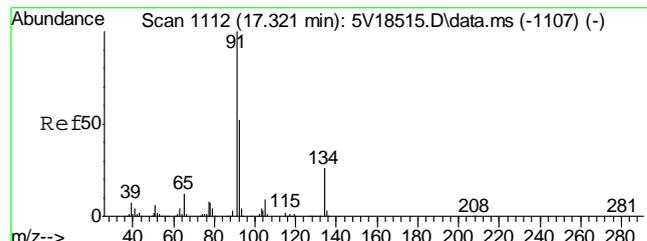
Tgt Ion	Resp	Lower	Upper
105	14122		
120	47.2	43.8	65.8



#86  
 p-Isopropyltoluene  
 Concen: 0.70 ug/l  
 RT: 16.944 min Scan# 1079  
 Delta R.T. -0.000 min  
 Lab File: 5V22147.D  
 Acq: 25 Jun 2012 7:38 pm

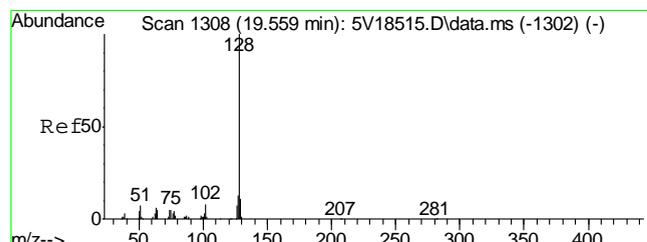
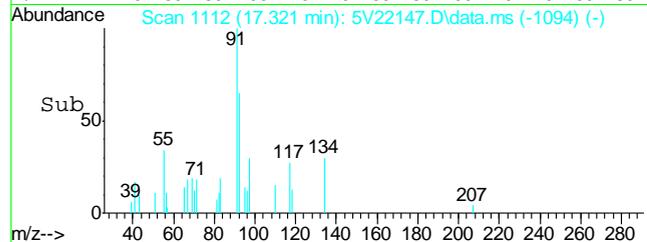
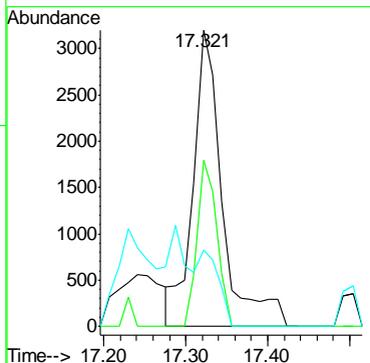
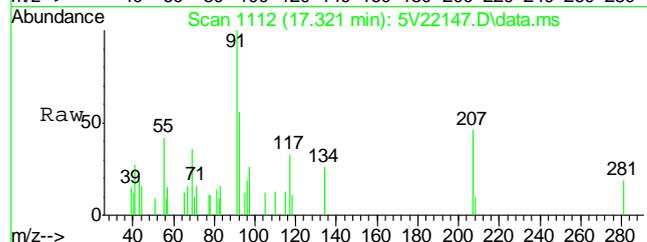
Tgt Ion	Resp	Lower	Upper
119	13177		
134	26.2	21.3	31.9
91	27.5	19.0	28.6





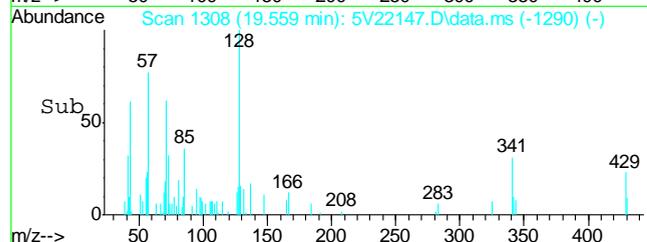
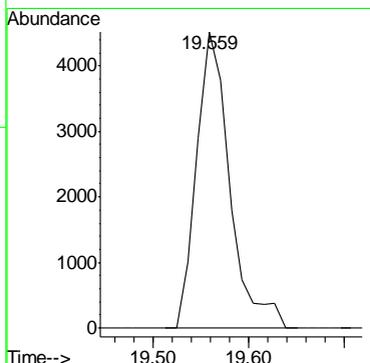
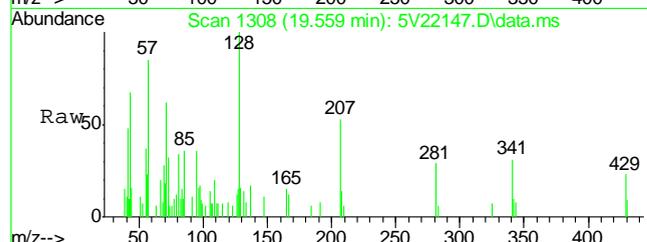
#88  
 n-Butylbenzene  
 Concen: 0.46 ug/l  
 RT: 17.321 min Scan# 1112  
 Delta R.T. -0.000 min  
 Lab File: 5V22147.D  
 Acq: 25 Jun 2012 7:38 pm

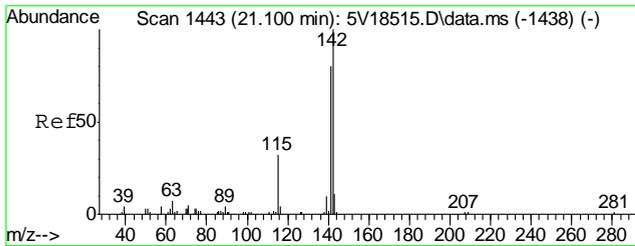
Tgt Ion	Resp	Lower	Upper
91	7925		
92	37.7	42.2	63.4#
134	42.7	21.4	32.2#



#91  
 Naphthalene  
 Concen: 1.40 ug/l  
 RT: 19.559 min Scan# 1308  
 Delta R.T. 0.001 min  
 Lab File: 5V22147.D  
 Acq: 25 Jun 2012 7:38 pm

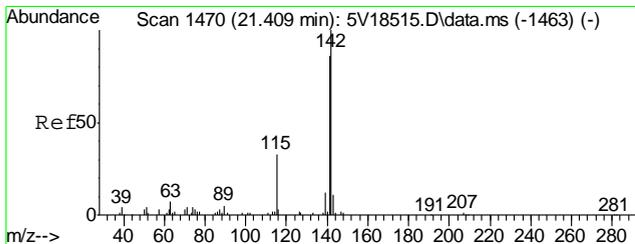
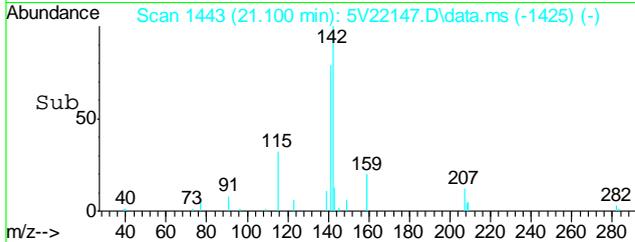
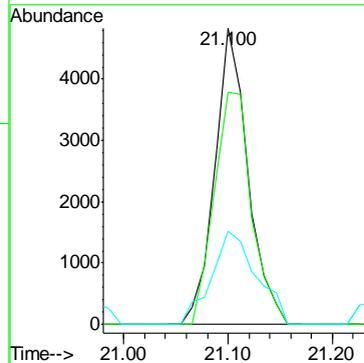
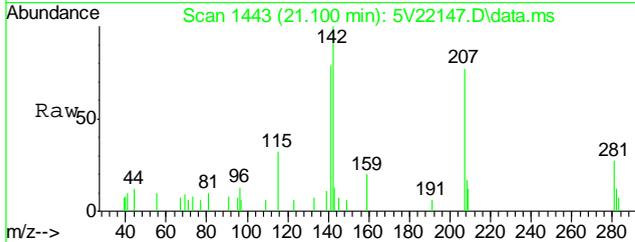
Tgt Ion	Resp
128	10872





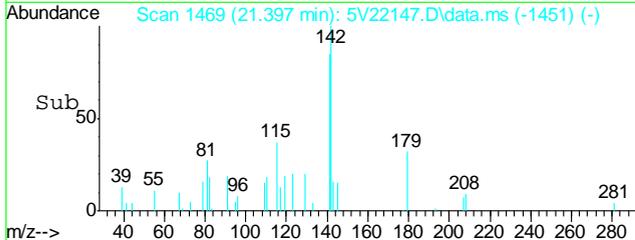
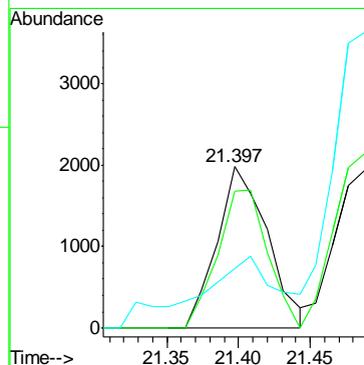
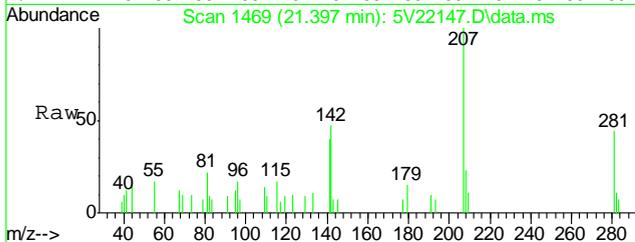
#94  
2-Methylnaphthalene  
Concen: 3.31 ug/l  
RT: 21.100 min Scan# 1443  
Delta R.T. 0.000 min  
Lab File: 5V22147.D  
Acq: 25 Jun 2012 7:38 pm

Tgt Ion	Resp	Lower	Upper
142	10686		
141	88.9	66.2	99.4
115	42.9	25.9	38.9#



#95  
1-Methylnaphthalene  
Concen: 2.09 ug/l  
RT: 21.397 min Scan# 1469  
Delta R.T. 0.000 min  
Lab File: 5V22147.D  
Acq: 25 Jun 2012 7:38 pm

Tgt Ion	Resp	Lower	Upper
142	4842		
141	84.8	68.9	103.3
115	72.4	27.3	40.9#



Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\V5062512.S\  
 Data File : 5V22134.D  
 Acq On : 25 Jun 2012 12:48 pm  
 Operator : BRETD  
 Sample : MB  
 Misc : MS4172,V5V1355,5.00,,100,5,1  
 ALS Vial : 3 Sample Multiplier: 1

Quant Time: Jun 26 08:09:10 2012  
 Quant Method : C:\msdchem\1\METHODS\V5AP1304TVH1304.M  
 Quant Title : 8260  
 QLast Update : Thu May 24 07:55:17 2012  
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
2) Pentafluorobenzene	11.647	168	184781	50.00	ug/l	0.00
35) 1,4-Difluorobenzene	12.446	114	295222	50.00	ug/l	0.00
53) Chlorobenzene-d5	15.095	117	382975	50.00	ug/l	0.00
74) 1,4-Dichlorobenzene-d4	17.070	152	250095	50.00	ug/l	0.00

System Monitoring Compounds

33) 1,2-Dichloroethane-d4	12.035	102	30897	54.76	ug/l	0.00
Spiked Amount	50.000	Range 70 - 130	Recovery	=	109.52%	
61) Toluene-d8	13.850	98	570641	44.09	ug/l	0.00
Spiked Amount	50.000	Range 70 - 130	Recovery	=	88.18%	
69) 4-Bromofluorobenzene	16.042	95	234606	44.25	ug/l	0.00
Spiked Amount	50.000	Range 70 - 130	Recovery	=	88.50%	

Target Compounds

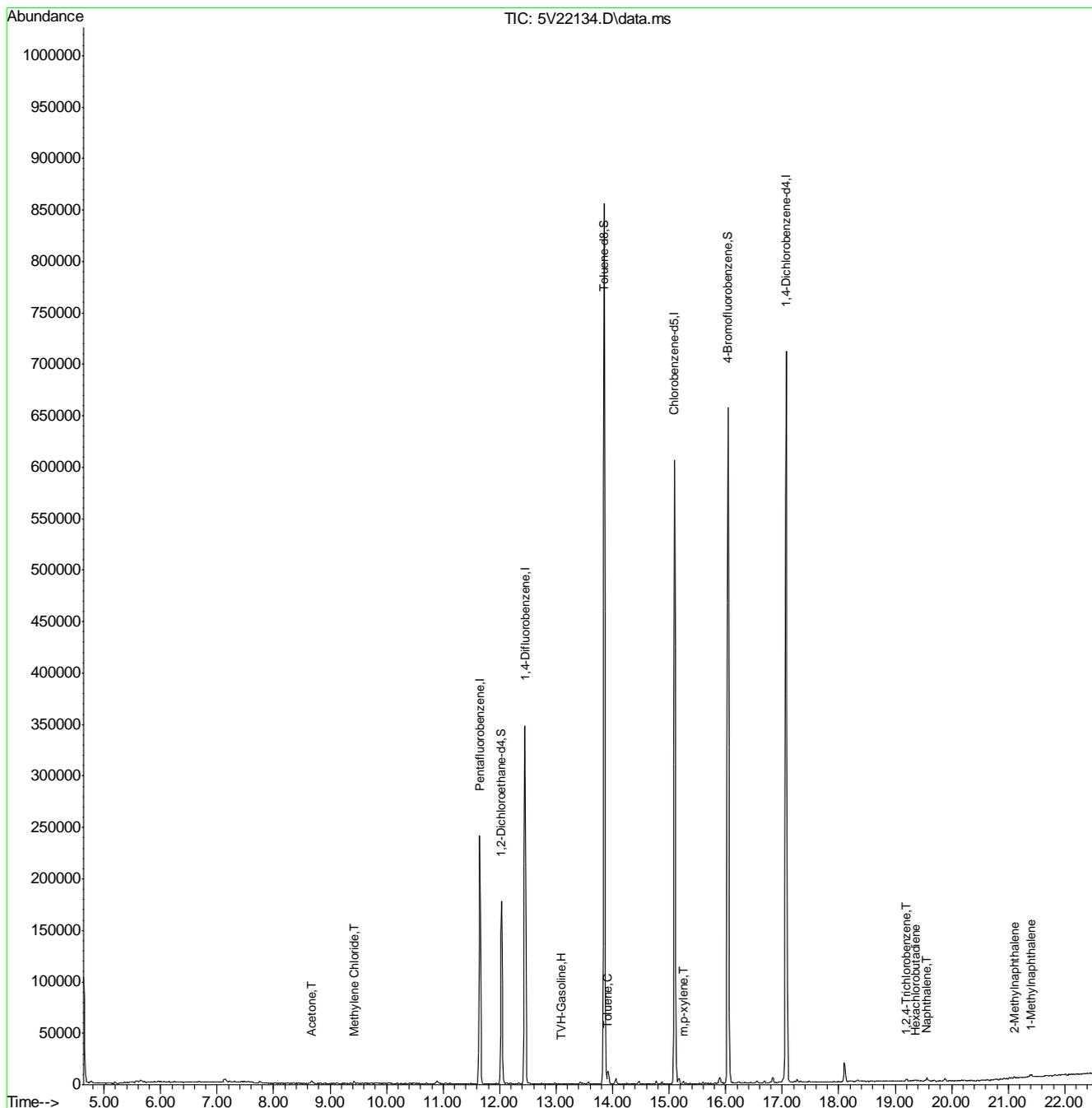
					Qvalue
1) TVH-Gasoline	13.102	TIC	3202m	1.56	ug/l
15) Acetone	8.679	58	1372	2.86	ug/l # 78
17) Methylene Chloride	9.421	84	1205	0.39	ug/l # 79
62) Toluene	13.908	92	2752	0.28	ug/l 95
72) m,p-xylene	15.255	106	1050	0.14	ug/l 96
90) 1,2,4-Trichlorobenzene	19.194	180	1591	0.24	ug/l # 86
91) Naphthalene	19.559	128	4770	0.96	ug/l 100
92) Hexachlorobutadiene	19.353	225	489	0.10	ug/l # 19
94) 2-Methylnaphthalene	21.112	142	1809	1.69	ug/l # 89
95) 1-Methylnaphthalene	21.397	142	1948	1.57	ug/l 94

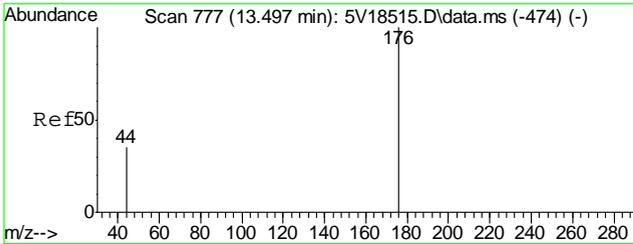
(#) = qualifier out of range (m) = manual integration (+) = signals summed

Quantitation Report (QT Reviewed)

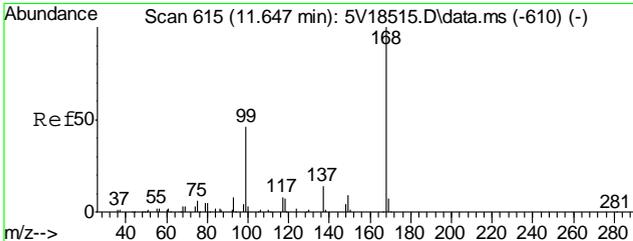
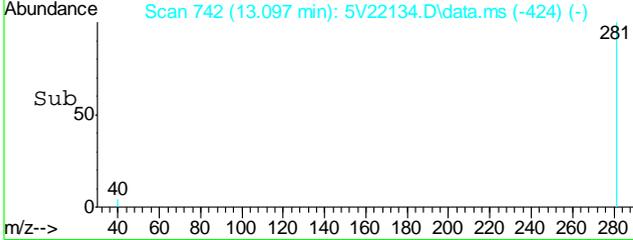
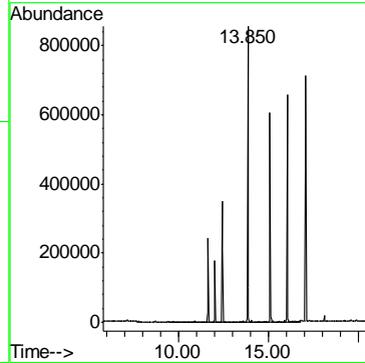
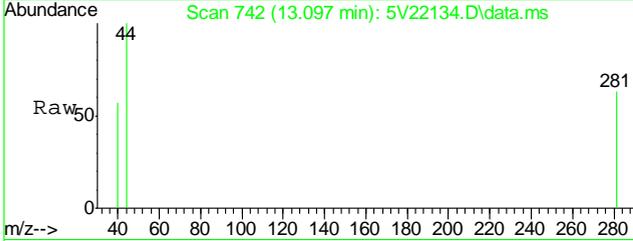
Data Path : C:\msdchem\1\DATA\V5062512.S\  
 Data File : 5V22134.D  
 Acq On : 25 Jun 2012 12:48 pm  
 Operator : BRETD  
 Sample : MB  
 Misc : MS4172,V5V1355,5.00,,100,5,1  
 ALS Vial : 3 Sample Multiplier: 1

Quant Time: Jun 26 08:09:10 2012  
 Quant Method : C:\msdchem\1\METHODS\V5AP1304TVH1304.M  
 Quant Title : 8260  
 QLast Update : Thu May 24 07:55:17 2012  
 Response via : Initial Calibration

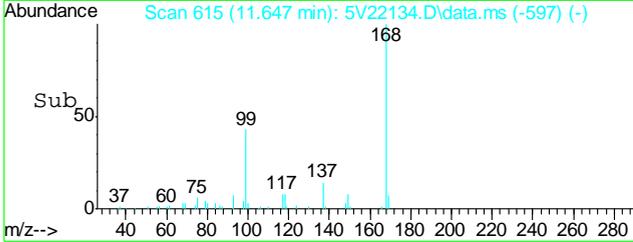
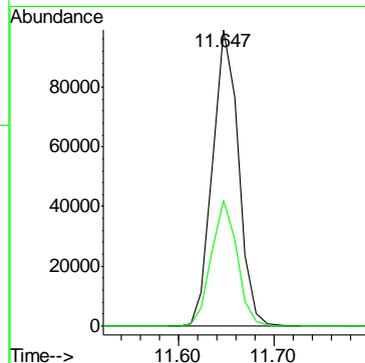
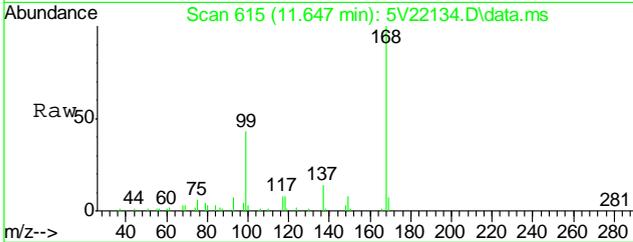


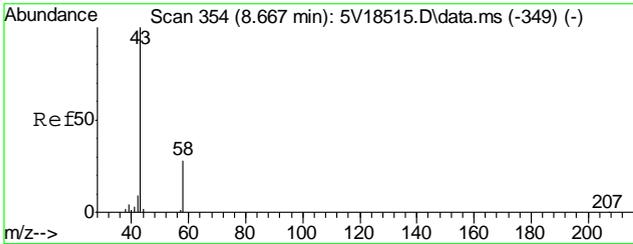


#1  
TVH-Gasoline  
Concen: 1.56 ug/l m  
RT: 13.102 min Scan# 742  
Delta R.T. 0.000 min  
Lab File: 5V22134.D  
Acq: 25 Jun 2012 12:48 pm  
Tgt Ion:TIC Resp: 3202



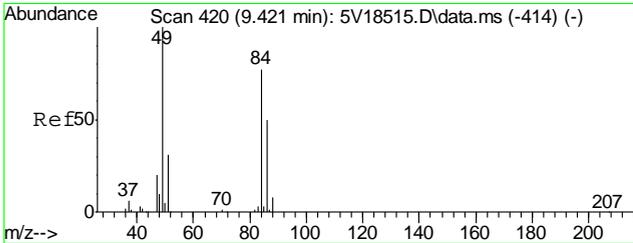
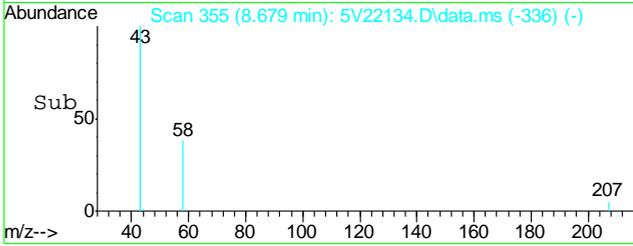
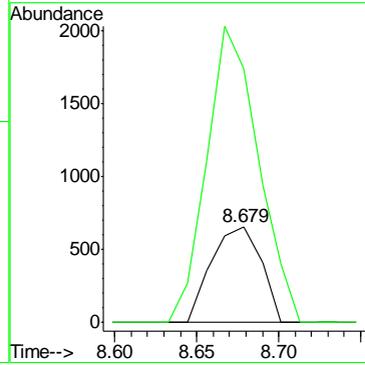
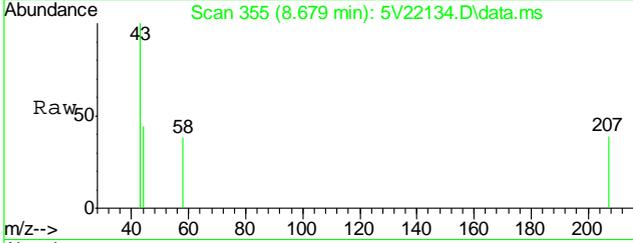
#2  
Pentafluorobenzene  
Concen: 50.00 ug/l  
RT: 11.647 min Scan# 615  
Delta R.T. -0.000 min  
Lab File: 5V22134.D  
Acq: 25 Jun 2012 12:48 pm  
Tgt Ion:168 Resp: 184781  
Ion Ratio Lower Upper  
168 100  
99 42.0 37.4 56.2





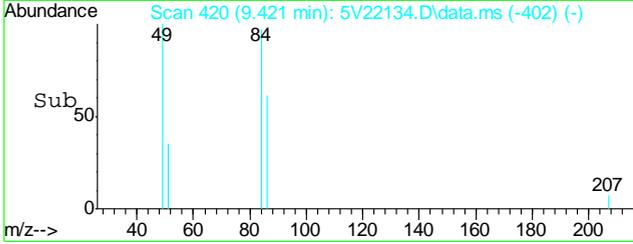
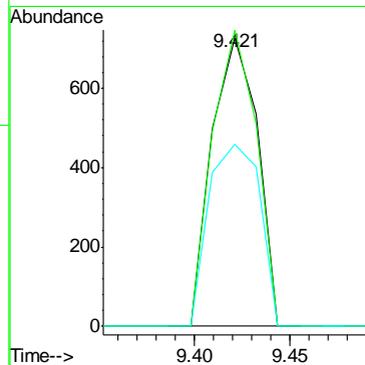
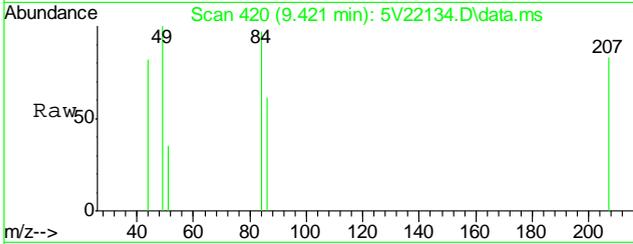
#15  
Acetone  
Concen: 2.86 ug/l  
RT: 8.679 min Scan# 355  
Delta R.T. 0.012 min  
Lab File: 5V22134.D  
Acq: 25 Jun 2012 12:48 pm

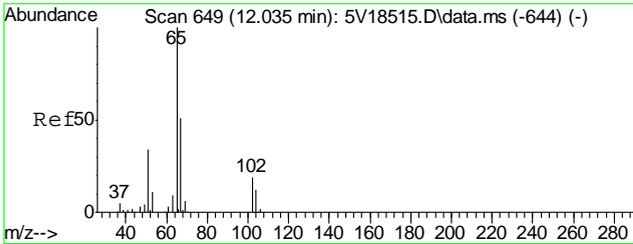
Tgt Ion: 58 Resp: 1372  
Ion Ratio Lower Upper  
58 100  
43 323.3 353.6 393.6#



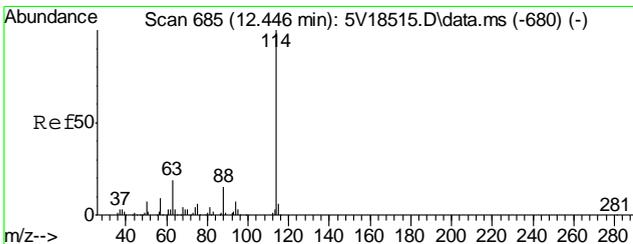
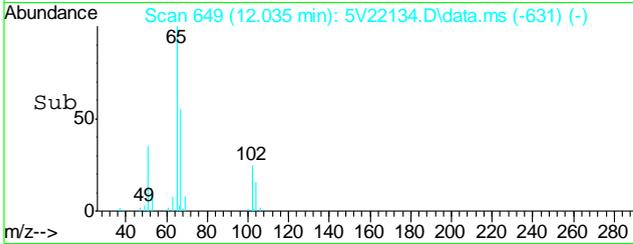
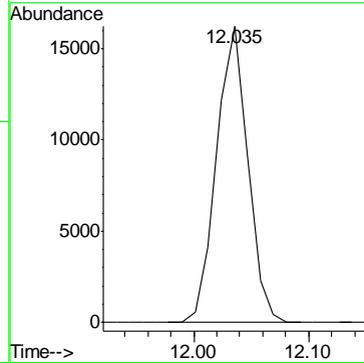
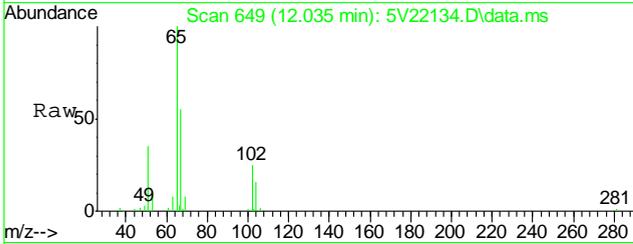
#17  
Methylene Chloride  
Concen: 0.39 ug/l  
RT: 9.421 min Scan# 420  
Delta R.T. -0.000 min  
Lab File: 5V22134.D  
Acq: 25 Jun 2012 12:48 pm

Tgt Ion: 84 Resp: 1205  
Ion Ratio Lower Upper  
84 100  
49 99.4 110.4 150.4#  
86 71.0 44.0 84.0

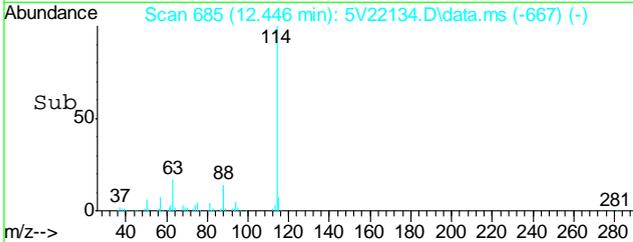
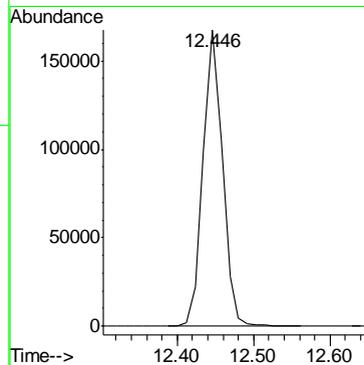
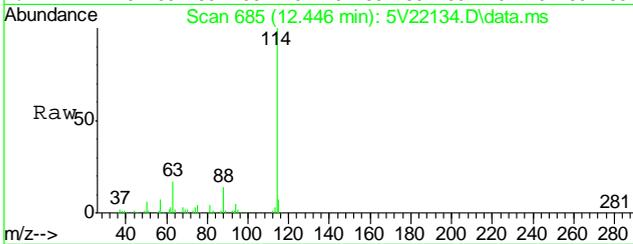


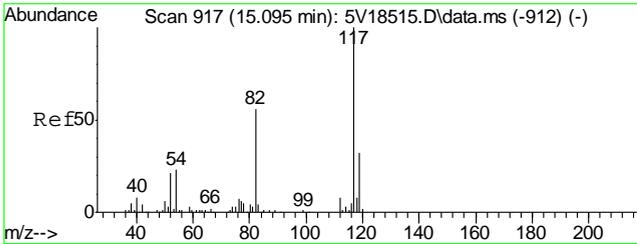


#33  
 1,2-Dichloroethane-d4  
 Concen: 54.76 ug/l  
 RT: 12.035 min Scan# 649  
 Delta R.T. 0.000 min  
 Lab File: 5V22134.D  
 Acq: 25 Jun 2012 12:48 pm  
 Tgt Ion:102 Resp: 30897

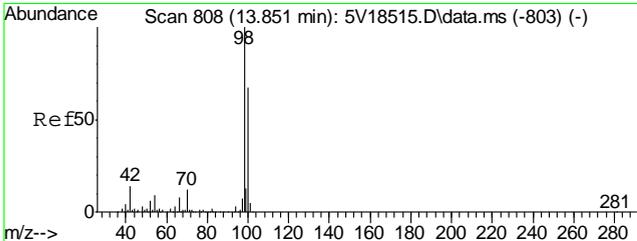
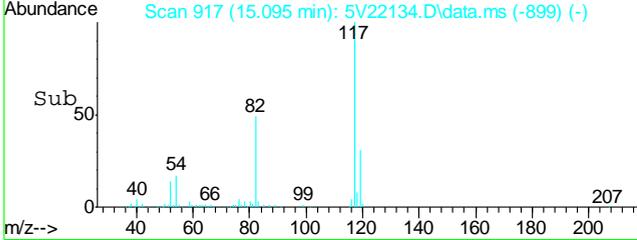
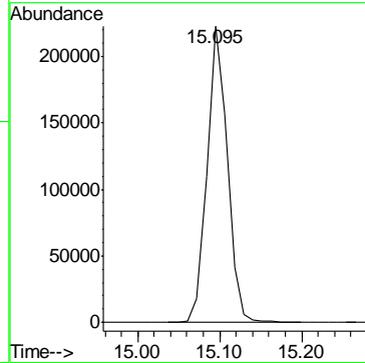
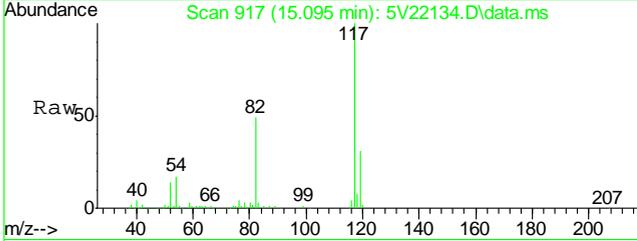


#35  
 1,4-Difluorobenzene  
 Concen: 50.00 ug/l  
 RT: 12.446 min Scan# 685  
 Delta R.T. -0.000 min  
 Lab File: 5V22134.D  
 Acq: 25 Jun 2012 12:48 pm  
 Tgt Ion:114 Resp: 295222

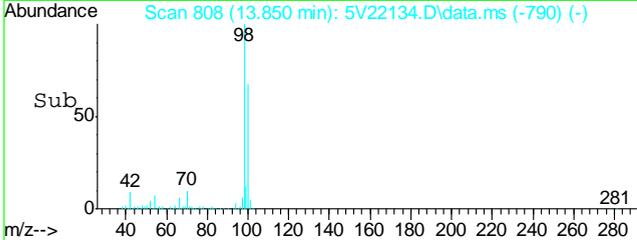
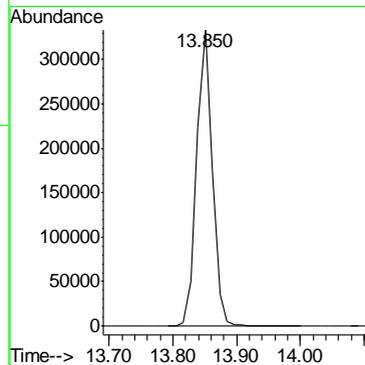
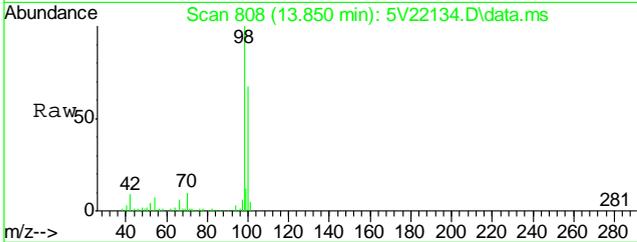


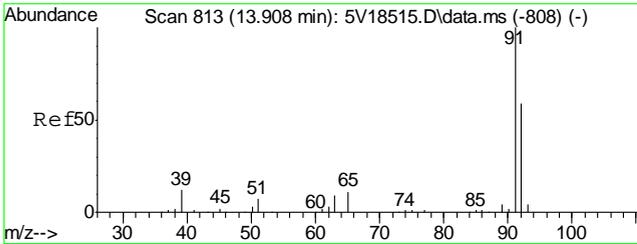


#53  
Chlorobenzene-d5  
Concen: 50.00 ug/l  
RT: 15.095 min Scan# 917  
Delta R.T. -0.000 min  
Lab File: 5V22134.D  
Acq: 25 Jun 2012 12:48 pm  
Tgt Ion:117 Resp: 382975



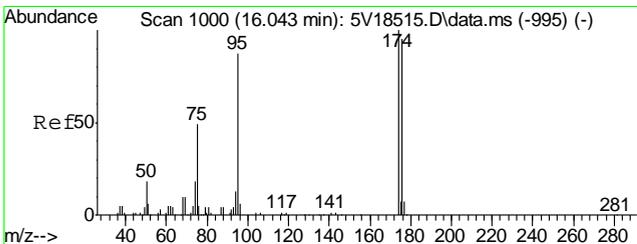
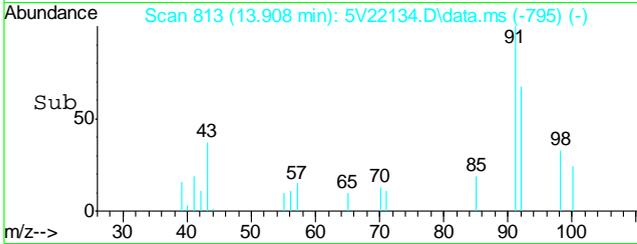
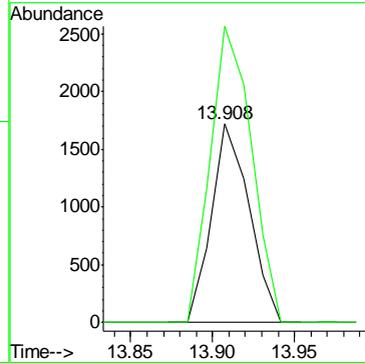
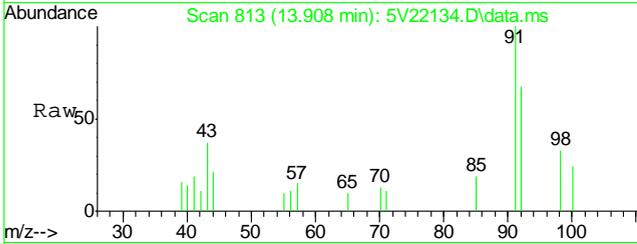
#61  
Toluene-d8  
Concen: 44.09 ug/l  
RT: 13.850 min Scan# 808  
Delta R.T. -0.000 min  
Lab File: 5V22134.D  
Acq: 25 Jun 2012 12:48 pm  
Tgt Ion: 98 Resp: 570641





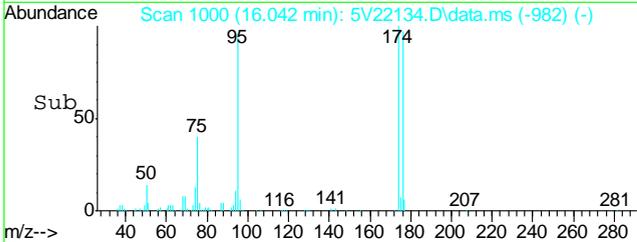
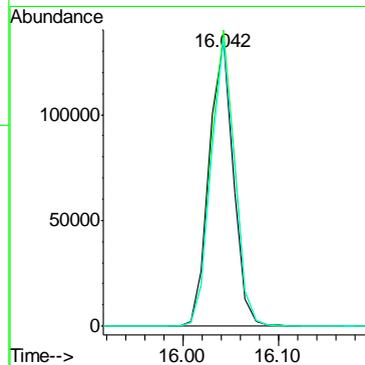
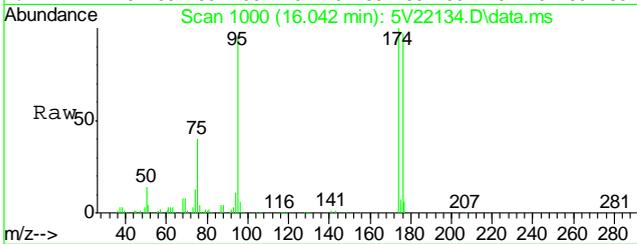
#62  
Toluene  
Concen: 0.28 ug/l  
RT: 13.908 min Scan# 813  
Delta R.T. -0.000 min  
Lab File: 5V22134.D  
Acq: 25 Jun 2012 12:48 pm

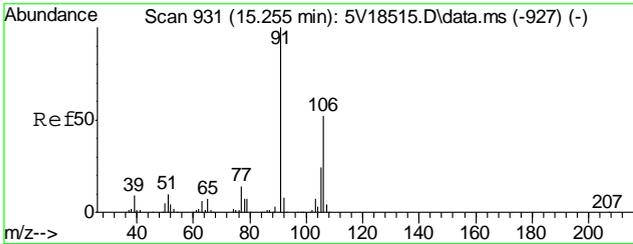
Tgt Ion: 92 Resp: 2752  
Ion Ratio Lower Upper  
92 100  
91 163.1 149.8 189.8



#69  
4-Bromofluorobenzene  
Concen: 44.25 ug/l  
RT: 16.042 min Scan# 1000  
Delta R.T. -0.000 min  
Lab File: 5V22134.D  
Acq: 25 Jun 2012 12:48 pm

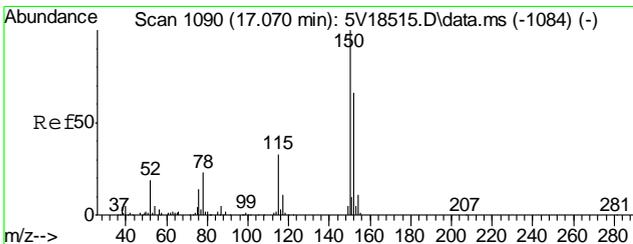
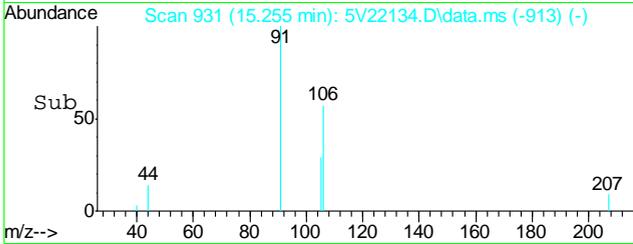
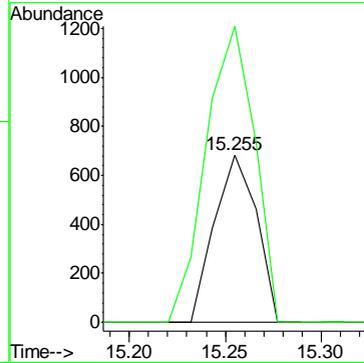
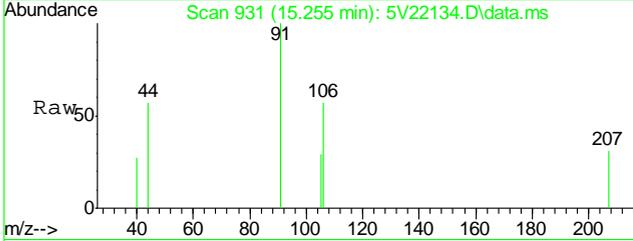
Tgt Ion: 95 Resp: 234606  
Ion Ratio Lower Upper  
95 100  
174 102.3 77.1 117.1  
176 98.2 73.4 113.4





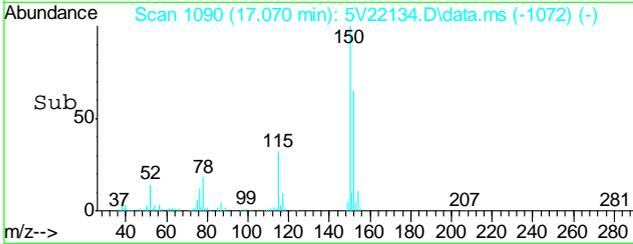
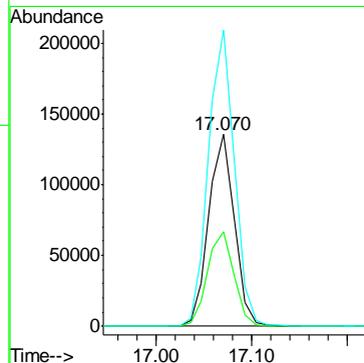
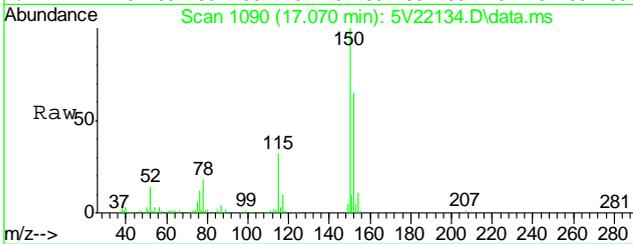
#72  
 m,p-xylene  
 Concen: 0.14 ug/l  
 RT: 15.255 min Scan# 931  
 Delta R.T. -0.000 min  
 Lab File: 5V22134.D  
 Acq: 25 Jun 2012 12:48 pm

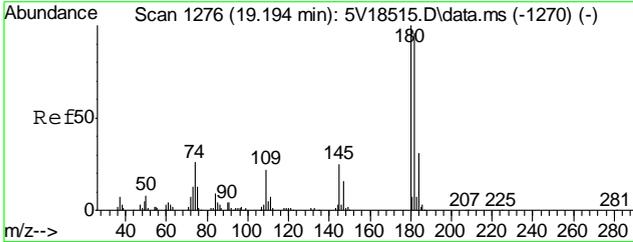
Tgt Ion	Resp	Lower	Upper
106	1050		
106	100		
91	203.7	177.1	217.1



#74  
 1,4-Dichlorobenzene-d4  
 Concen: 50.00 ug/l  
 RT: 17.070 min Scan# 1090  
 Delta R.T. -0.000 min  
 Lab File: 5V22134.D  
 Acq: 25 Jun 2012 12:48 pm

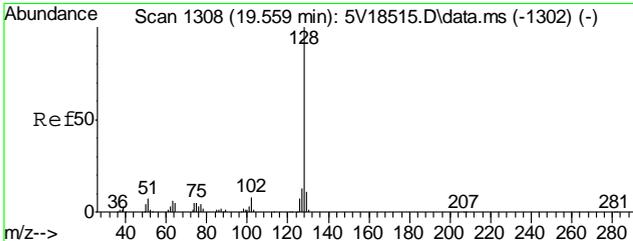
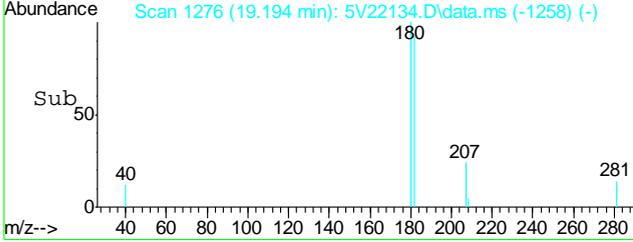
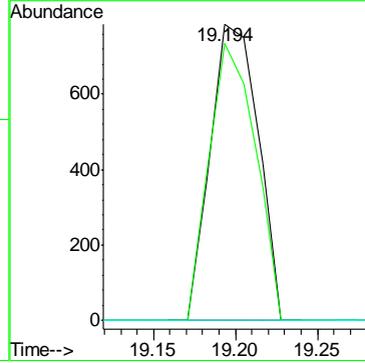
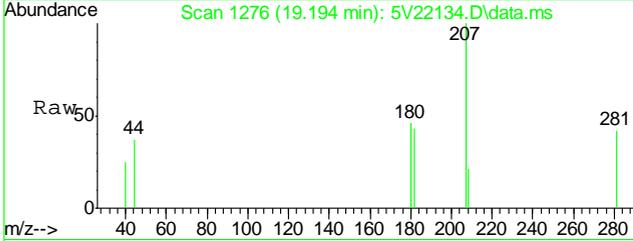
Tgt Ion	Resp	Lower	Upper
152	250095		
152	100		
115	50.7	41.4	62.0
150	156.7	153.9	230.9





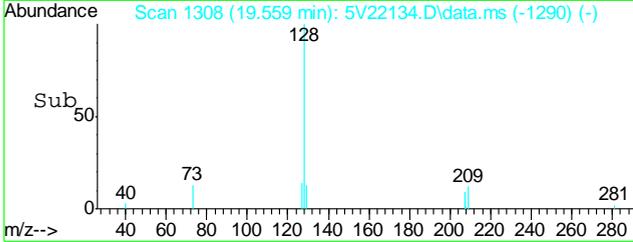
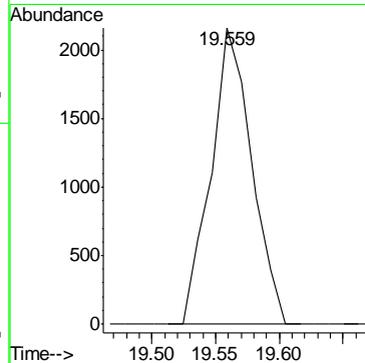
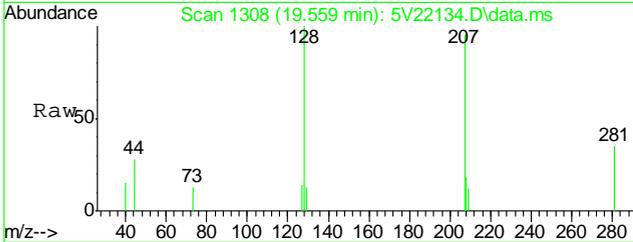
#90  
 1,2,4-Trichlorobenzene  
 Concen: 0.24 ug/l  
 RT: 19.194 min Scan# 1276  
 Delta R.T. -0.000 min  
 Lab File: 5V22134.D  
 Acq: 25 Jun 2012 12:48 pm

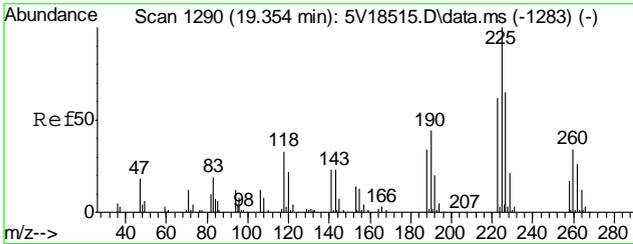
Tgt Ion	Resp	Lower	Upper
180	1591	100	
182	90.6	76.2	114.4
145	0.0	20.1	30.1#



#91  
 Naphthalene  
 Concen: 0.96 ug/l  
 RT: 19.559 min Scan# 1308  
 Delta R.T. 0.001 min  
 Lab File: 5V22134.D  
 Acq: 25 Jun 2012 12:48 pm

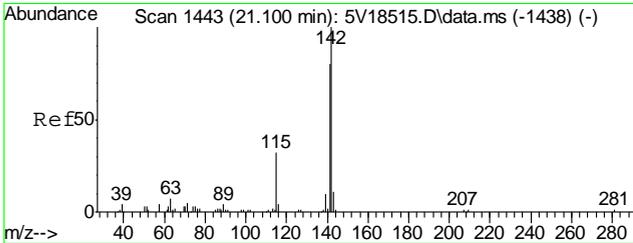
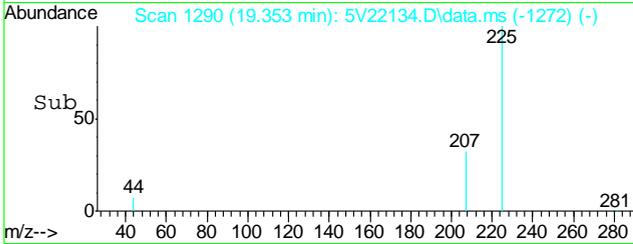
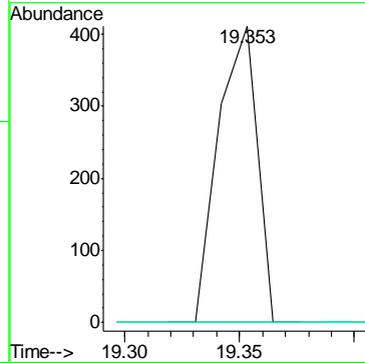
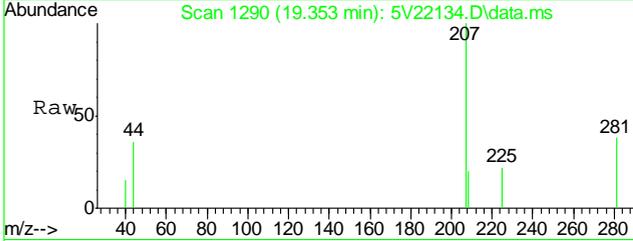
Tgt Ion	Resp
128	4770





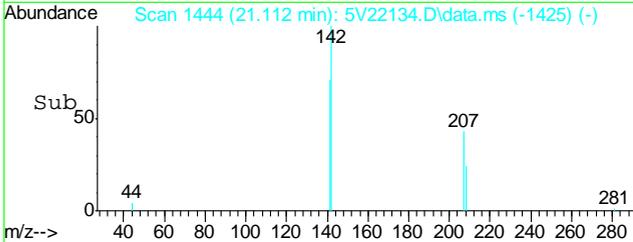
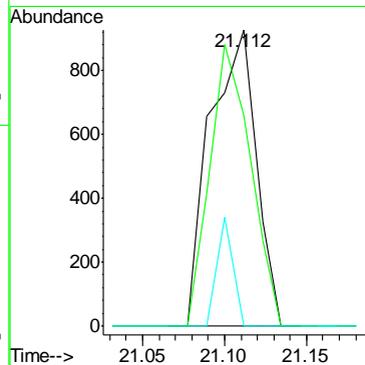
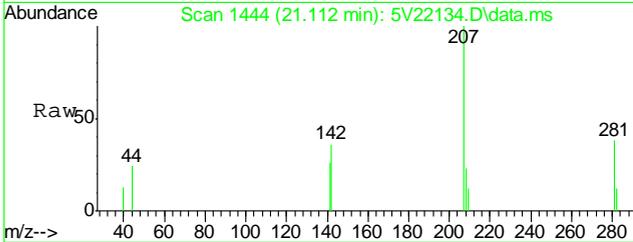
#92  
Hexachlorobutadiene  
Concen: 0.10 ug/l  
RT: 19.353 min Scan# 1290  
Delta R.T. -0.000 min  
Lab File: 5V22134.D  
Acq: 25 Jun 2012 12:48 pm

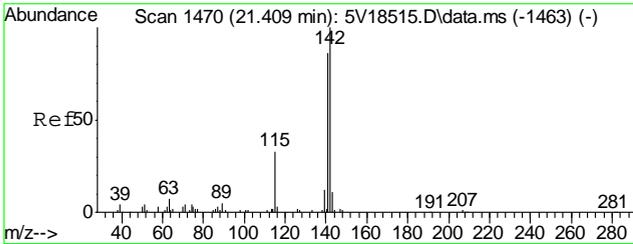
Tgt Ion	Resp	Lower	Upper
225	489		
225	100		
223	0.0	49.5	74.3#
227	0.0	51.1	76.7#



#94  
2-Methylnaphthalene  
Concen: 1.69 ug/l  
RT: 21.112 min Scan# 1444  
Delta R.T. 0.012 min  
Lab File: 5V22134.D  
Acq: 25 Jun 2012 12:48 pm

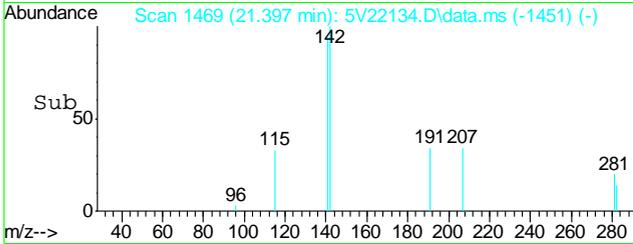
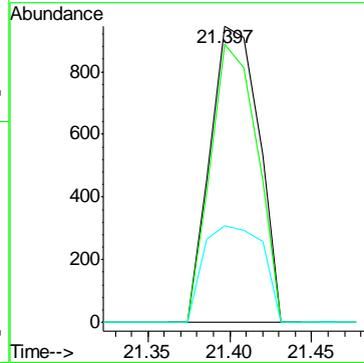
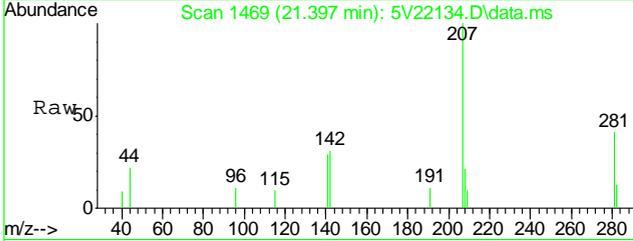
Tgt Ion	Resp	Lower	Upper
142	1809		
142	100		
141	84.2	66.2	99.4
115	12.9	25.9	38.9#





#95  
 1-Methylnaphthalene  
 Concen: 1.57 ug/l  
 RT: 21.397 min Scan# 1469  
 Delta R.T. 0.000 min  
 Lab File: 5V22134.D  
 Acq: 25 Jun 2012 12:48 pm

Tgt Ion	Resp	Lower	Upper
142	100		
141	90.2	68.9	103.3
115	39.6	27.3	40.9



6.2.1  
6

## 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:** D35710  
**Account:** XTOKRWR XTO Energy  
**Project:** FRU 297-8B

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
OP6113-MB	3G09818.D	1	06/22/12	SM	06/22/12	OP6113	E3G436

The QC reported here applies to the following samples:

Method: SW846 8270C BY SIM

D35710-1

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

CAS No.	Surrogate Recoveries	Limits
4165-60-0	Nitrobenzene-d5	76% 10-145%
321-60-8	2-Fluorobiphenyl	80% 10-130%
1718-51-0	Terphenyl-d14	97% 22-130%

# Blank Spike Summary

**Job Number:** D35710  
**Account:** XTOKRWR XTO Energy  
**Project:** FRU 297-8B

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
OP6113-BS	3G09819.D	1	06/22/12	SM	06/22/12	OP6113	E3G436

The QC reported here applies to the following samples:

Method: SW846 8270C BY SIM

D35710-1

CAS No.	Compound	Spike ug/kg	BSP ug/kg	BSP %	Limits
83-32-9	Acenaphthene	83.3	75.2	90	34-130
120-12-7	Anthracene	83.3	81.4	98	35-130
56-55-3	Benzo(a)anthracene	83.3	78.2	94	36-130
50-32-8	Benzo(a)pyrene	83.3	65.7	79	36-130
205-99-2	Benzo(b)fluoranthene	83.3	65.2	78	35-130
207-08-9	Benzo(k)fluoranthene	83.3	90.2	108	37-130
218-01-9	Chrysene	83.3	76.5	92	40-130
53-70-3	Dibenzo(a,h)anthracene	83.3	56.5	68	32-130
206-44-0	Fluoranthene	83.3	78.6	94	38-130
86-73-7	Fluorene	83.3	76.8	92	35-130
193-39-5	Indeno(1,2,3-cd)pyrene	83.3	55.7	67	28-130
91-20-3	Naphthalene	83.3	75.2	90	35-130
129-00-0	Pyrene	83.3	76.5	92	37-130

CAS No.	Surrogate Recoveries	BSP	Limits
4165-60-0	Nitrobenzene-d5	79%	10-145%
321-60-8	2-Fluorobiphenyl	82%	10-130%
1718-51-0	Terphenyl-d14	87%	22-130%

\* = Outside of Control Limits.

7.2.1  
 7

# Matrix Spike/Matrix Spike Duplicate Summary

**Job Number:** D35710  
**Account:** XTOKRWR XTO Energy  
**Project:** FRU 297-8B

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
OP6113-MS	3G09821.D	1	06/22/12	SM	06/22/12	OP6113	E3G436
OP6113-MSD	3G09822.D	1	06/22/12	SM	06/22/12	OP6113	E3G436
D35708-1	3G09820.D	1	06/22/12	SM	06/22/12	OP6113	E3G436

The QC reported here applies to the following samples:

Method: SW846 8270C BY SIM

D35710-1

CAS No.	Compound	D35708-1 ug/kg	Spike Q	ug/kg	MS ug/kg	MS %	MSD ug/kg	MSD %	RPD	Limits Rec/RPD
83-32-9	Acenaphthene	ND	87.7	77.5	88	78.3	89	1	10-155/30	
120-12-7	Anthracene	ND	87.7	85.8	98	85.9	98	0	10-155/30	
56-55-3	Benzo(a)anthracene	ND	87.7	97.0	111	98.1	112	1	10-175/30	
50-32-8	Benzo(a)pyrene	ND	87.7	77.3	88	79.5	91	3	10-164/30	
205-99-2	Benzo(b)fluoranthene	ND	87.7	87.5	100	92.6	106	6	10-165/30	
207-08-9	Benzo(k)fluoranthene	ND	87.7	66.9	76	69.3	79	4	10-178/30	
218-01-9	Chrysene	ND	87.7	85.0	97	87.0	99	2	10-147/30	
53-70-3	Dibenzo(a,h)anthracene	ND	87.7	75.7	86	72.8	83	4	10-144/30	
206-44-0	Fluoranthene	ND	87.7	72.9	83	71.7	82	2	10-207/30	
86-73-7	Fluorene	ND	87.7	83.6	95	82.6	94	1	10-163/30	
193-39-5	Indeno(1,2,3-cd)pyrene	ND	87.7	71.2	81	70.3	80	1	10-180/30	
91-20-3	Naphthalene	57.8	87.7	130	82	129	81	1	10-198/30	
129-00-0	Pyrene	ND	87.7	105	120	97.5	111	7	10-189/30	

CAS No.	Surrogate Recoveries	MS	MSD	D35708-1	Limits
4165-60-0	Nitrobenzene-d5	60%	57%	62%	10-145%
321-60-8	2-Fluorobiphenyl	75%	72%	71%	10-130%
1718-51-0	Terphenyl-d14	89%	78%	90%	22-130%

\* = Outside of Control Limits.

GC/MS Semi-volatiles

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Raw Data

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## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\062212\  
 Data File : 3g09824.D  
 Acq On : 22 Jun 2012 8:03 pm  
 Operator : SARAHM1  
 Sample : D35710-1  
 Misc : OP6113,E3G436,30.15,,,1,1  
 ALS Vial : 17 Sample Multiplier: 1

Quant Time: Jun 25 10:28:15 2012  
 Quant Method : C:\msdchem\1\METHODS\SIMPE3G436.M  
 Quant Title : PAHSIM BASE  
 QLast Update : Fri Jun 22 16:07:04 2012  
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Naphthalene-d8	6.429	136	247896	4.0000	ug/mL	0.00
6) Acenaphthene-d10	8.823	164	142669	4.0000	ug/mL	0.00
14) Phenanthrene-d10	11.369	188	183193	4.0000	ug/mL	0.00
18) Chrysene-d12	16.423	240	140159	4.0000	ug/mL	0.00
23) Perylene-d12	18.997	264	120082	4.0000	ug/mL	0.00

## System Monitoring Compounds

2) Nitrobenzene-d5	5.718	82	1117196	31.4093	ug/mL	-0.01
Spiked Amount	50.000	Range 25 - 135	Recovery =	62.82%		
7) 2-Fluorobiphenyl	7.807	172	1845380	38.9641	ug/mL	-0.01
Spiked Amount	50.000	Range 25 - 135	Recovery =	77.92%		
20) Terphenyl-d14	14.472	244	1071623	42.1726	ug/mL	-0.02
Spiked Amount	50.000	Range 25 - 135	Recovery =	84.34%		

## Target Compounds

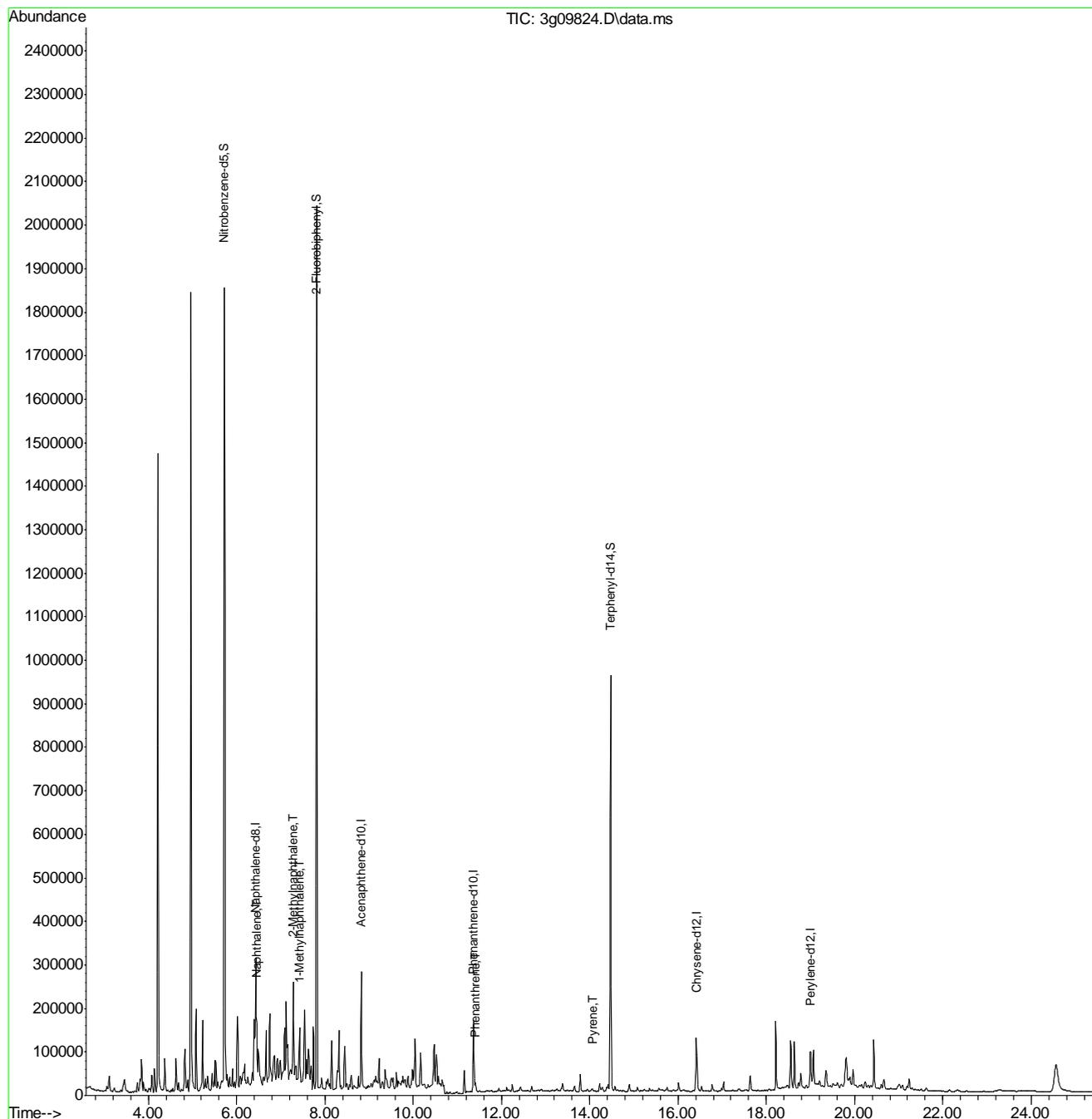
					Qvalue
3) N-Nitrosodimethylamine	0.000	74	0	N.D.	d
4) N-Nitrosodi-propylamine	0.000	70	0	N.D.	d
5) Naphthalene	6.454	128	103686	1.5246	ug/mL 95
8) 2-Methylnaphthalene	7.276	142	129373	3.1776	ug/mL 97
9) 1-Methylnaphthalene	7.426	142	62003	1.5433	ug/mL 93
10) Acenaphthylene	0.000	152	0	N.D.	d
11) Acenaphthene	0.000	154	0	N.D.	d
12) Fluorene	0.000	166	0	N.D.	d
13) Diphenylamine	0.000	169	0	N.D.	d
15) Phenanthrene	11.409	178	24548	0.4344	ug/mL 94
16) Anthracene	0.000	178	0	N.D.	d
17) Fluoranthene	0.000	202	0	N.D.	d
19) Pyrene	14.060	202	4513	0.0885	ug/mL 87
21) Benzo(a)anthracene	0.000	228	0	N.D.	d
22) Chrysene	0.000	228	0	N.D.	d
24) Benzo(b)fluoranthene	0.000	252	0	N.D.	d
25) Benzo(k)fluoranthene	0.000	252	0	N.D.	d
26) Benzo(a)pyrene	0.000	252	0	N.D.	d
27) Indeno(1,2,3-cd)pyrene	0.000	276	0	N.D.	d
28) Dibenzo(a,h)anthracene	0.000	278	0	N.D.	d
29) Benzo(g,h,i)perylene	0.000	276	0	N.D.	d

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

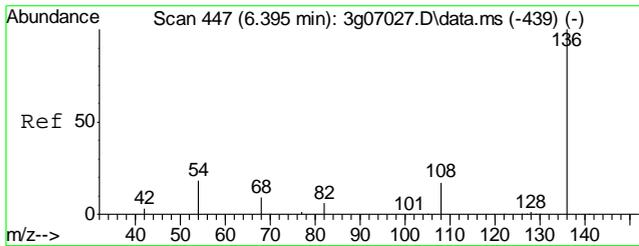
Quantitation Report (QT Reviewed)

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 Operator : SARAHM1  
 Sample : D35710-1  
 Misc : OP6113,E3G436,30.15,,,1,1  
 ALS Vial : 17 Sample Multiplier: 1

Quant Time: Jun 25 10:28:15 2012  
 Quant Method : C:\msdchem\1\METHODS\SIMPE3G436.M  
 Quant Title : PAHSIM BASE  
 QLast Update : Fri Jun 22 16:07:04 2012  
 Response via : Initial Calibration

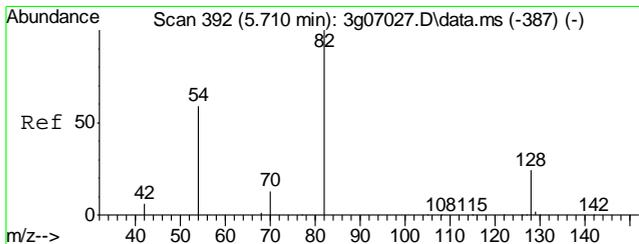
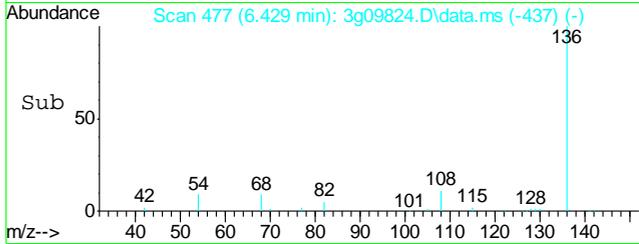
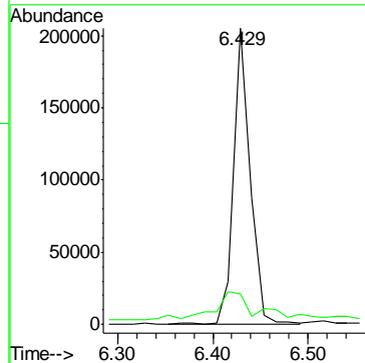
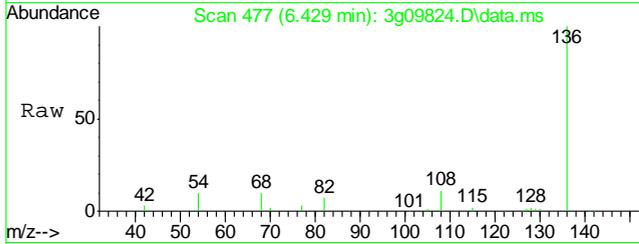


8.1.1  
8



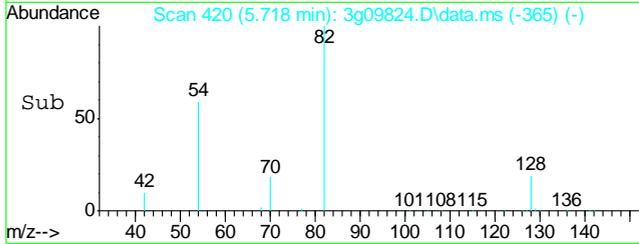
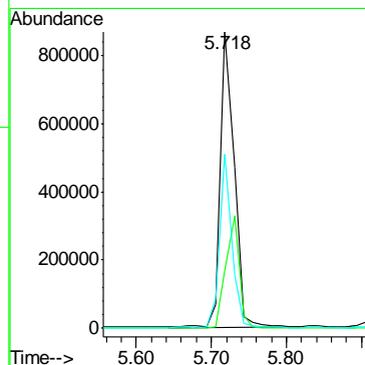
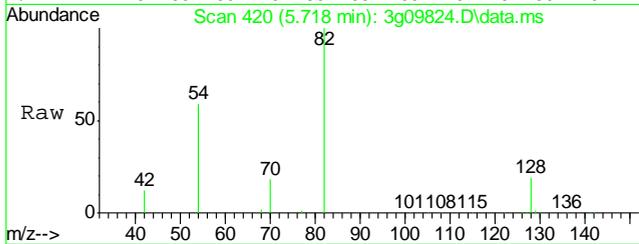
#1  
 Naphthalene-d8  
 Concen: 4.0000 ug/mL  
 RT: 6.429 min Scan# 477  
 Delta R.T. 0.000 min  
 Lab File: 3g09824.D  
 Acq: 22 Jun 12 8:03 pm

Tgt Ion	Resp	Lower	Upper
136	247896	100	
68	28.0	0.0	32.0

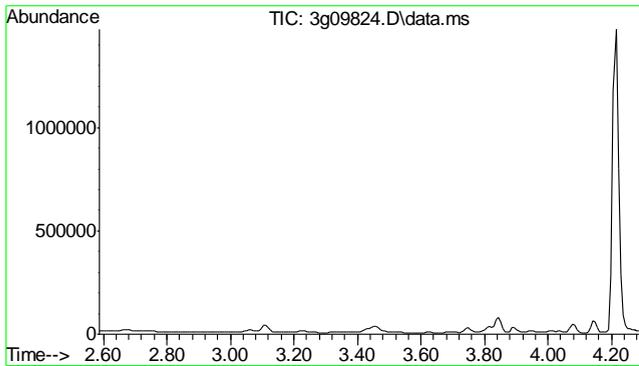


#2  
 Nitrobenzene-d5  
 Concen: 31.4093 ug/mL  
 RT: 5.718 min Scan# 420  
 Delta R.T. -0.012 min  
 Lab File: 3g09824.D  
 Acq: 22 Jun 12 8:03 pm

Tgt Ion	Resp	Lower	Upper
82	1117196	100	
128	36.7	15.2	55.2
54	52.0	34.0	74.0



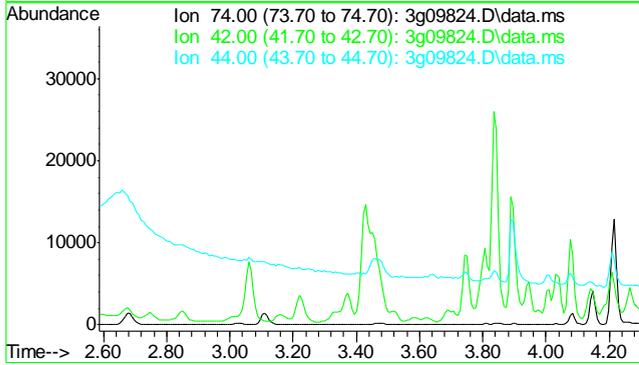
8.1.1  
 8



#3  
 N-Nitrosodimethylamine  
 Concen: N.D. ug/mL  
 Expected RT: 2.79 min

Lab File: 3g09824.D  
 Acq: 22 Jun 12 8:03 pm

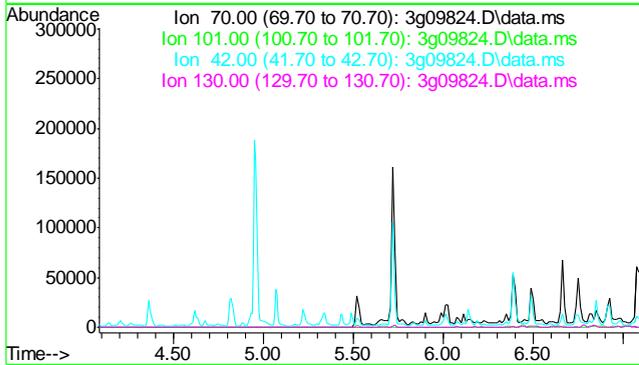
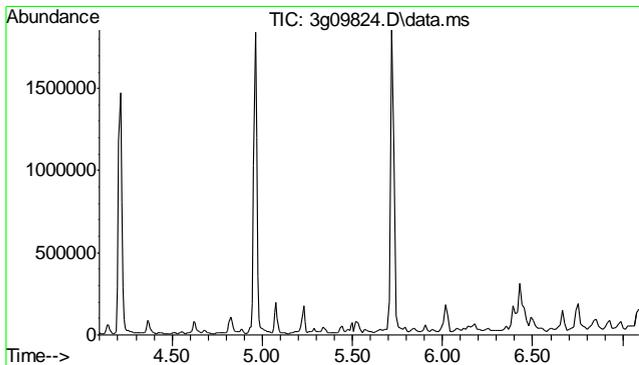
Tgt Ion:	74
Sig	Exp Ratio
74	100
42	51.8
44	10.0

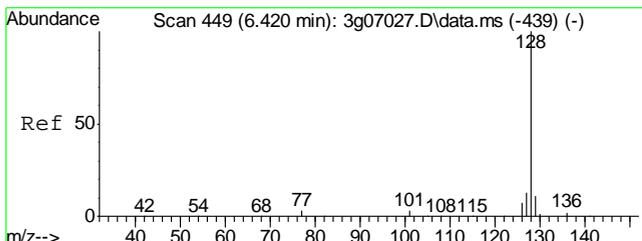


#4  
 N-Nitrosodi-propylamine  
 Concen: N.D. ug/mL  
 Expected RT: 5.59 min

Lab File: 3g09824.D  
 Acq: 22 Jun 12 8:03 pm

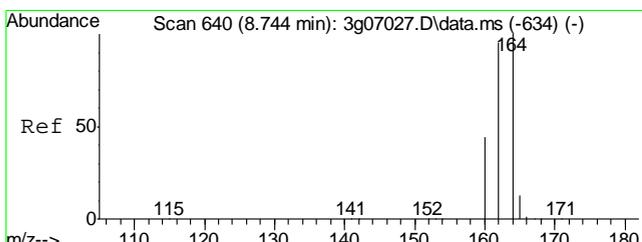
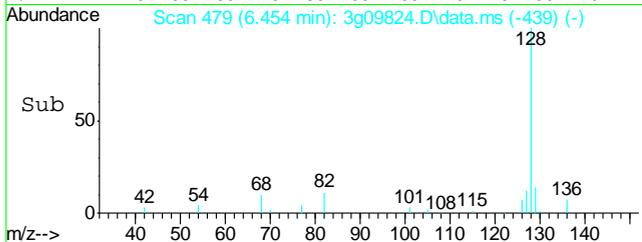
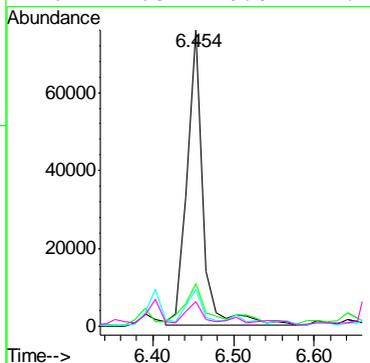
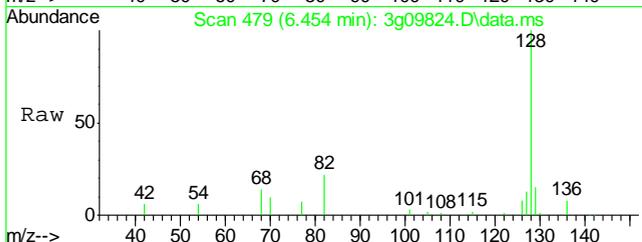
Tgt Ion:	70
Sig	Exp Ratio
70	100
101	10.3
42	59.5
130	20.1





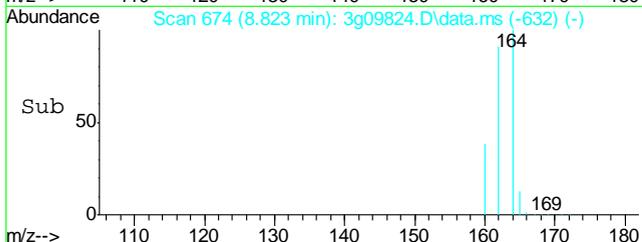
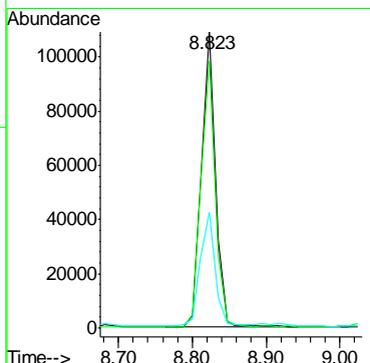
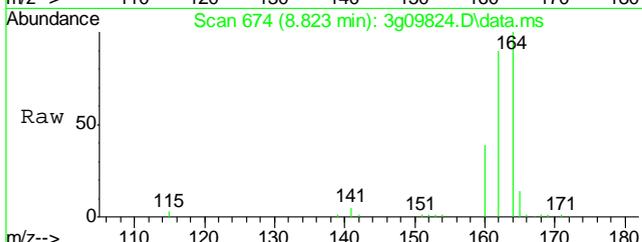
#5  
 Naphthalene  
 Concen: 1.5246 ug/mL  
 RT: 6.454 min Scan# 479  
 Delta R.T. 0.000 min  
 Lab File: 3g09824.D  
 Acq: 22 Jun 12 8:03 pm

Tgt Ion	Resp	Lower	Upper
128	103686		
129	15.4	0.0	30.9
127	12.7	0.0	32.4
126	7.3	0.0	27.7

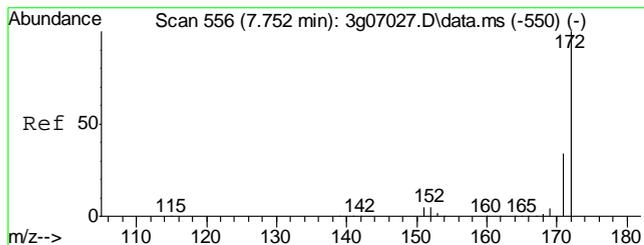


#6  
 Acenaphthene-d10  
 Concen: 4.0000 ug/mL  
 RT: 8.823 min Scan# 674  
 Delta R.T. 0.000 min  
 Lab File: 3g09824.D  
 Acq: 22 Jun 12 8:03 pm

Tgt Ion	Resp	Lower	Upper
164	142669		
162	91.9	72.5	112.5
160	41.7	20.9	60.9

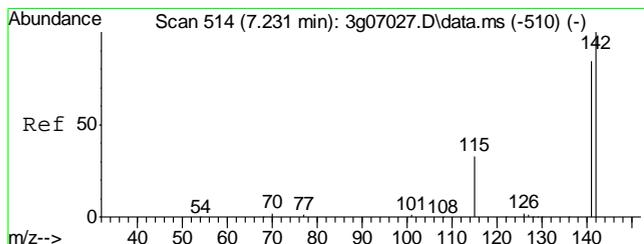
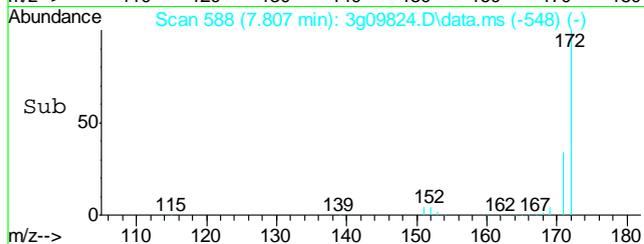
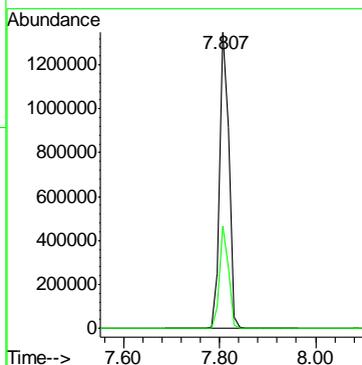
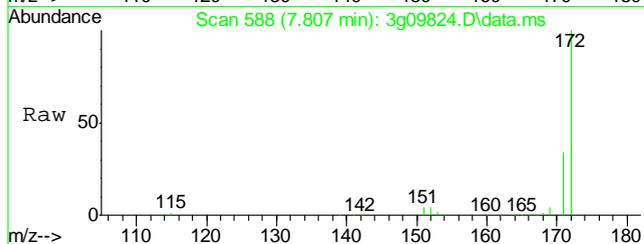


8.1.1  
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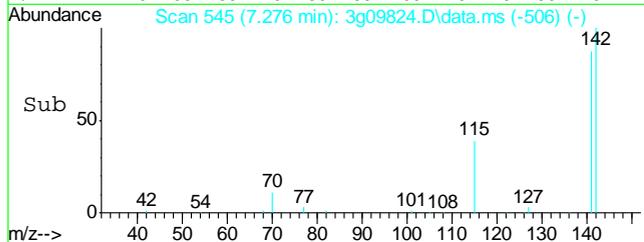
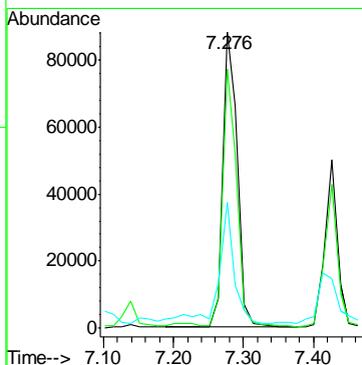
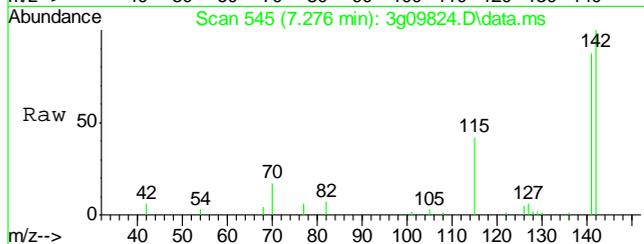
#7  
 2-Fluorobiphenyl  
 Concen: 38.9641 ug/mL  
 RT: 7.807 min Scan# 588  
 Delta R.T. -0.012 min  
 Lab File: 3g09824.D  
 Acq: 22 Jun 12 8:03 pm

Tgt Ion: 172 Resp: 1845380  
 Ion Ratio Lower Upper  
 172 100  
 171 33.1 13.2 53.2

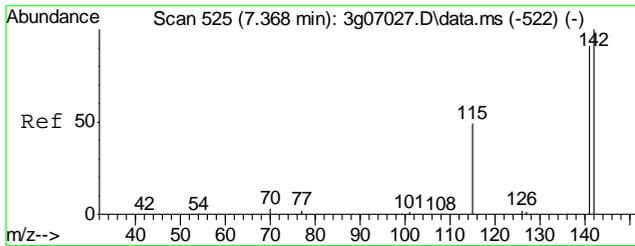


#8  
 2-Methylnaphthalene  
 Concen: 3.1776 ug/mL  
 RT: 7.276 min Scan# 545  
 Delta R.T. -0.012 min  
 Lab File: 3g09824.D  
 Acq: 22 Jun 12 8:03 pm

Tgt Ion: 142 Resp: 129373  
 Ion Ratio Lower Upper  
 142 100  
 141 83.5 63.1 103.1  
 115 39.4 14.8 54.8

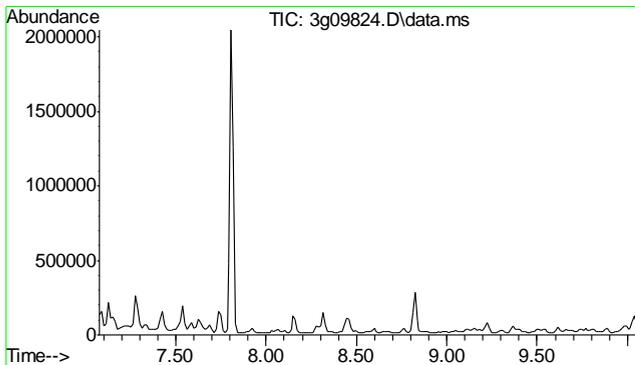
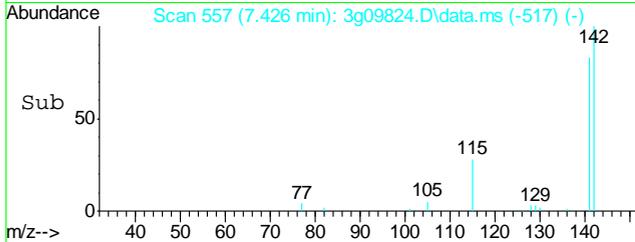
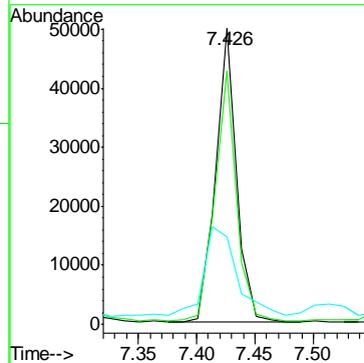
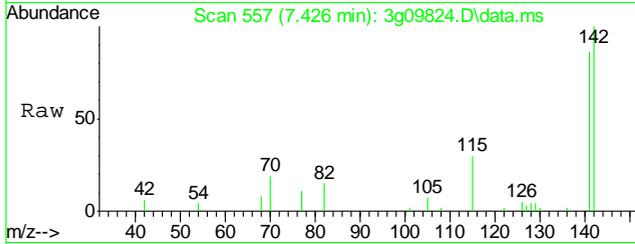


8.1.1  
 8



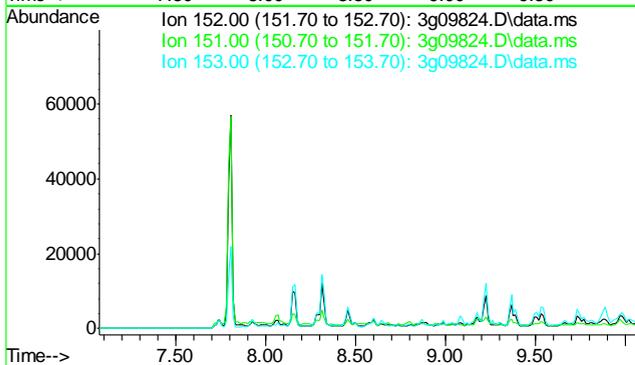
#9  
 1-Methylnaphthalene  
 Concen: 1.5433 ug/mL  
 RT: 7.426 min Scan# 557  
 Delta R.T. 0.000 min  
 Lab File: 3g09824.D  
 Acq: 22 Jun 12 8:03 pm

Tgt Ion	Resp	Lower	Upper
142	62003	100	
141	88.2	66.4	106.4
115	47.0	16.1	56.1

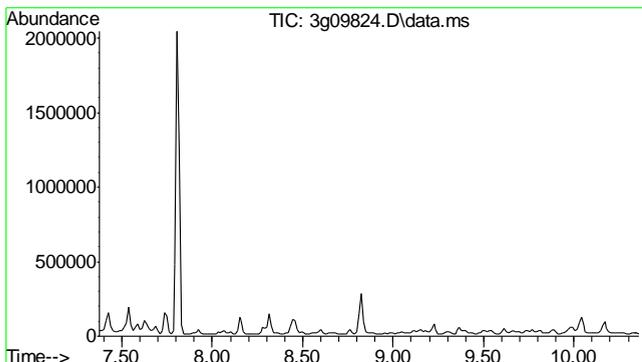


#10  
 Acenaphthylene  
 Concen: N.D. ug/mL  
 Expected RT: 8.58 min  
 Lab File: 3g09824.D  
 Acq: 22 Jun 12 8:03 pm

Tgt Ion	Sig	Exp Ratio
152	100	
151		19.1
153		14.1



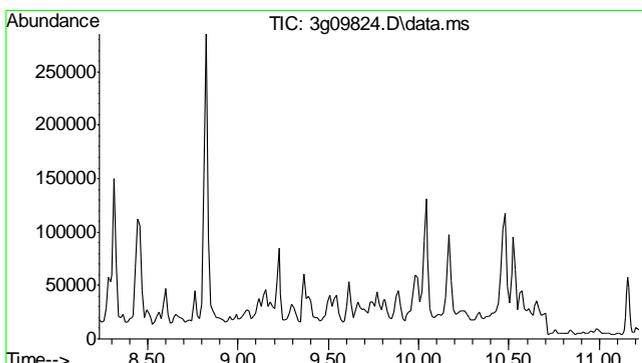
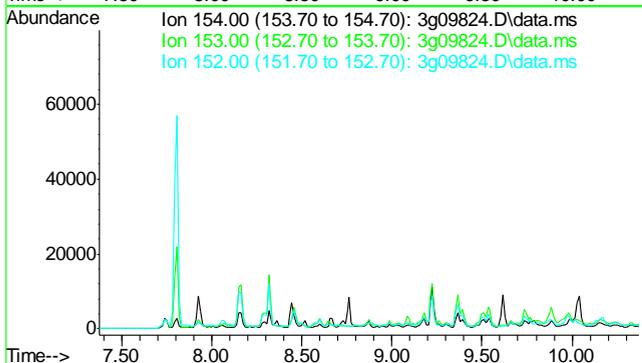
8.1.1  
 8



#11  
 Acenaphthene  
 Concen: N.D. ug/mL  
 Expected RT: 8.87 min

Lab File: 3g09824.D  
 Acq: 22 Jun 12 8:03 pm

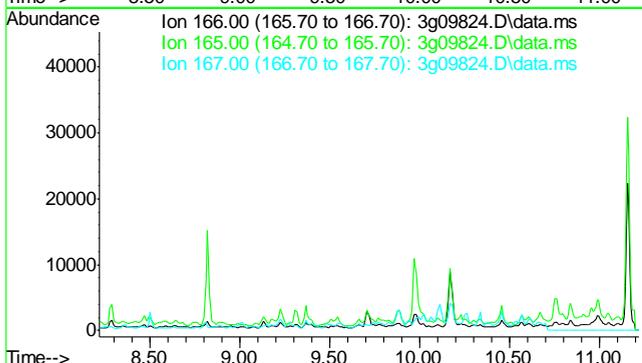
Tgt Ion	Exp Ratio
154	100
153	104.2
152	45.3



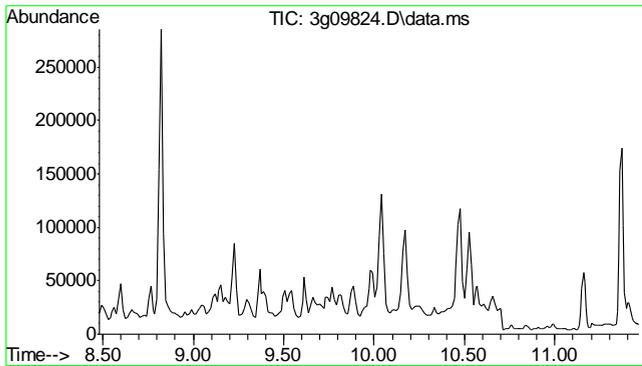
#12  
 Fluorene  
 Concen: N.D. ug/mL  
 Expected RT: 9.72 min

Lab File: 3g09824.D  
 Acq: 22 Jun 12 8:03 pm

Tgt Ion	Exp Ratio
166	100
165	90.8
167	13.3

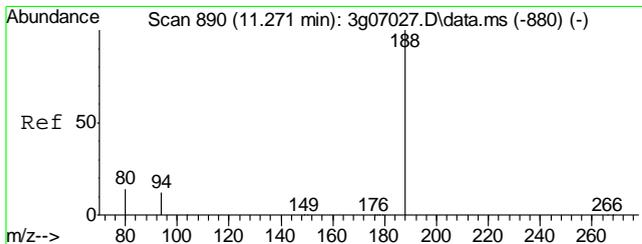
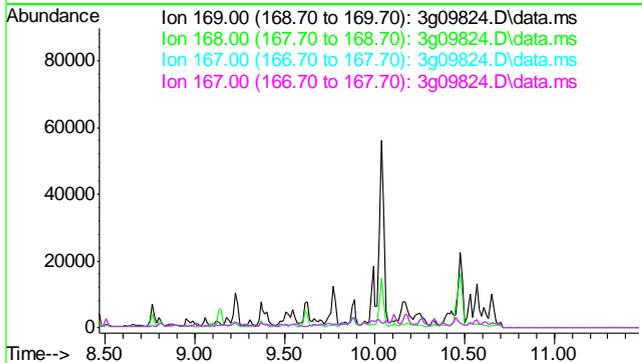


8.1.1  
8



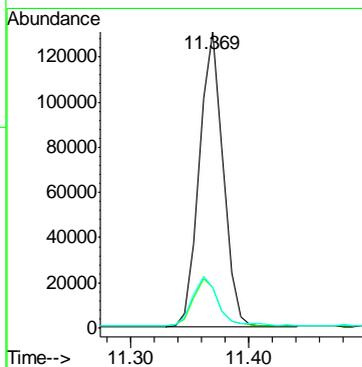
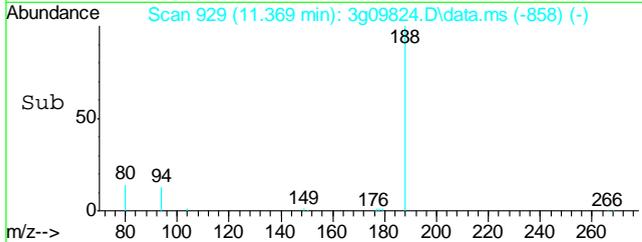
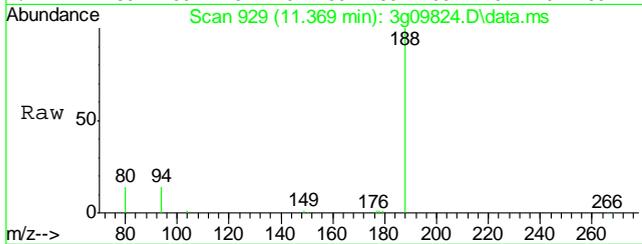
#13  
 Diphenylamine  
 Concen: N.D. ug/mL  
 Expected RT: 9.97 min  
 Lab File: 3g09824.D  
 Acq: 22 Jun 12 8:03 pm

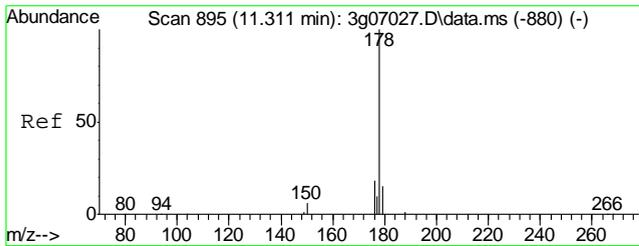
Tgt Ion	Exp Ratio
169	100
168	62.1
167	33.2
167	33.2



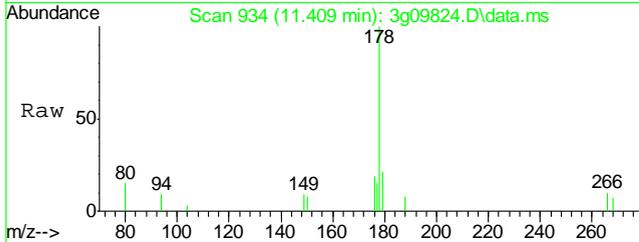
#14  
 Phenanthrene-d10  
 Concen: 4.0000 ug/mL  
 RT: 11.369 min Scan# 929  
 Delta R.T. 0.000 min  
 Lab File: 3g09824.D  
 Acq: 22 Jun 12 8:03 pm

Tgt Ion	Resp	Ion Ratio	Lower	Upper
188	183193	100		
94		16.6	0.0	37.2
80		18.3	0.0	38.3

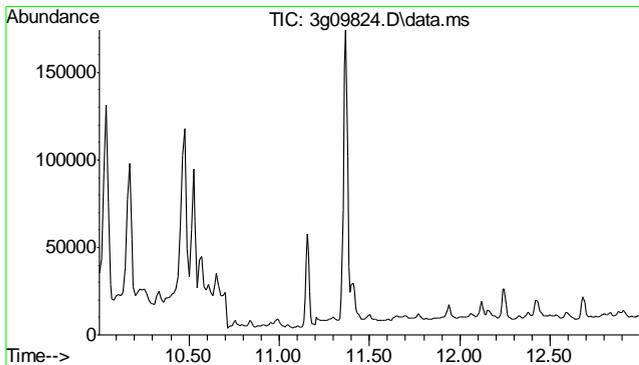
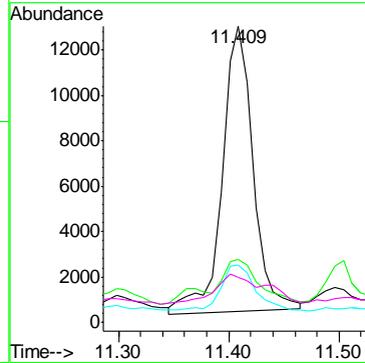
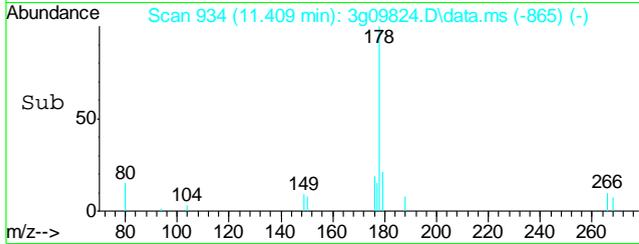




#15  
 Phenanthrene  
 Concen: 0.4344 ug/mL  
 RT: 11.409 min Scan# 934  
 Delta R.T. -0.008 min  
 Lab File: 3g09824.D  
 Acq: 22 Jun 12 8:03 pm

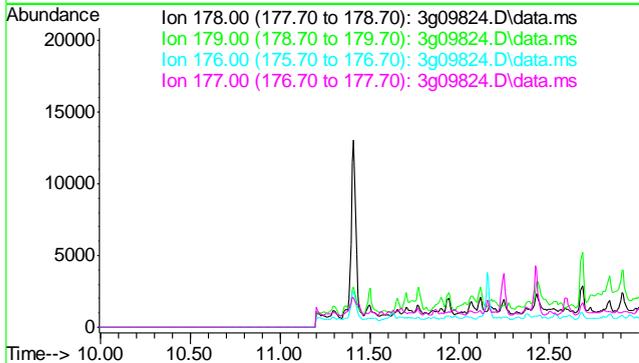


Tgt Ion	Resp	Lower	Upper
178	24548		
179	16.4	0.0	35.1
176	17.0	0.0	38.6
177	15.2	0.0	30.4

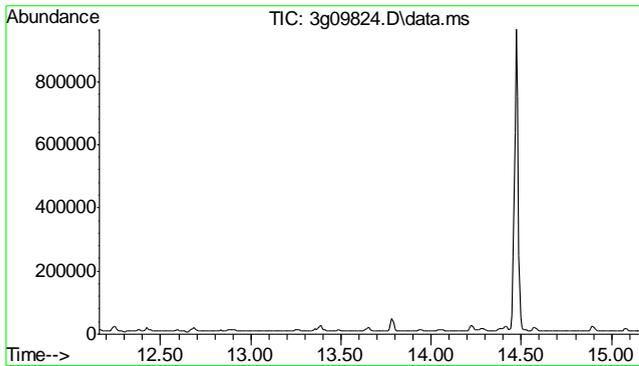


#16  
 Anthracene  
 Concen: N.D. ug/mL  
 Expected RT: 11.50 min  
 Lab File: 3g09824.D  
 Acq: 22 Jun 12 8:03 pm

Tgt Ion	Sig	Exp Ratio
178	100	
179	15.0	
176	17.9	
177	8.9	

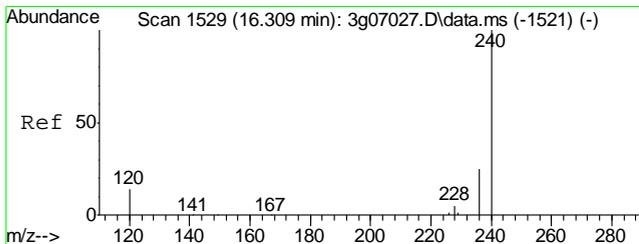
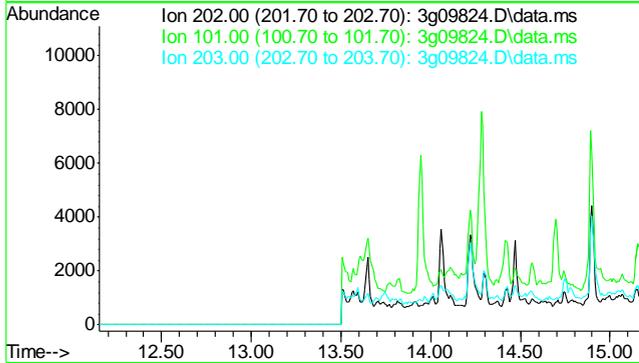


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#17  
 Fluoranthene  
 Concen: N.D. ug/mL  
 Expected RT: 13.66 min  
  
 Lab File: 3g09824.D  
 Acq: 22 Jun 12 8:03 pm  
  
 Tgt Ion: 202  

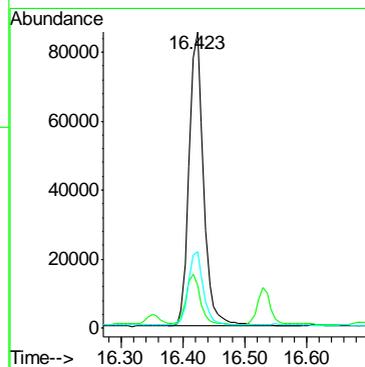
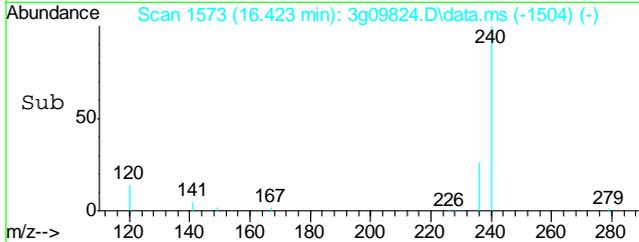
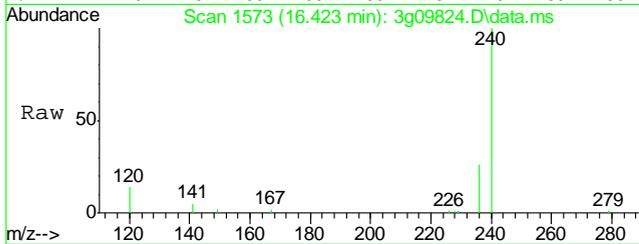
Sig	Exp Ratio
202	100
101	17.6
203	17.1

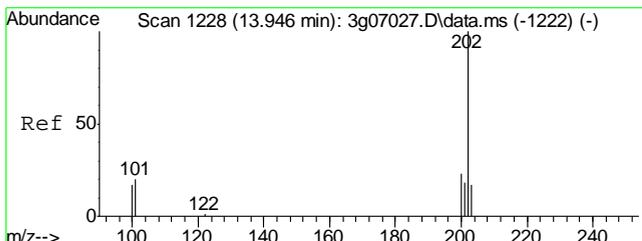


#18  
 Chrysene-d12  
 Concen: 4.0000 ug/mL  
 RT: 16.423 min Scan# 1573  
 Delta R.T. 0.000 min  
 Lab File: 3g09824.D  
 Acq: 22 Jun 12 8:03 pm

Tgt Ion: 240 Resp: 140159  

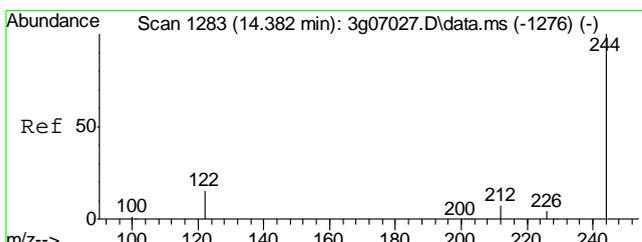
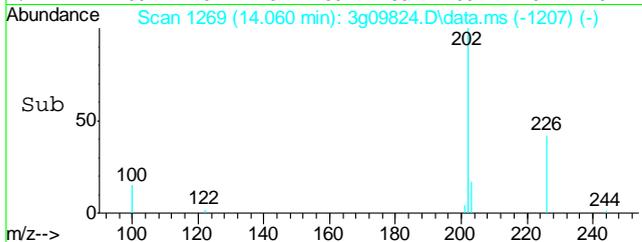
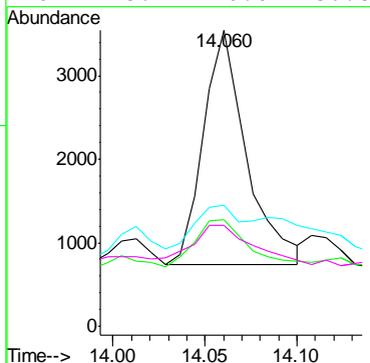
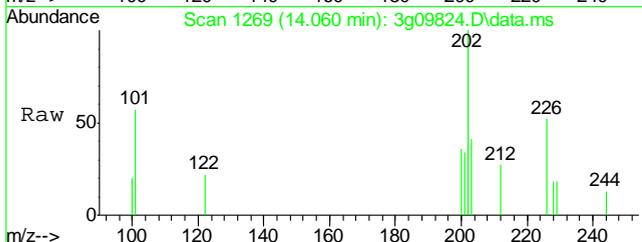
Ion	Ratio	Lower	Upper
240	100		
120	17.4	0.0	37.2
236	25.1	5.2	45.2





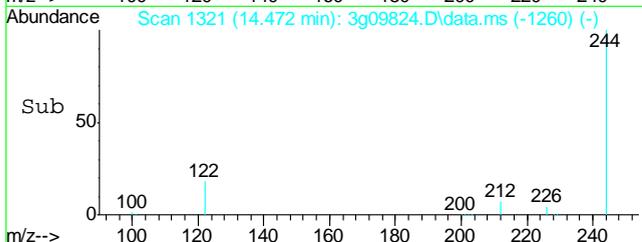
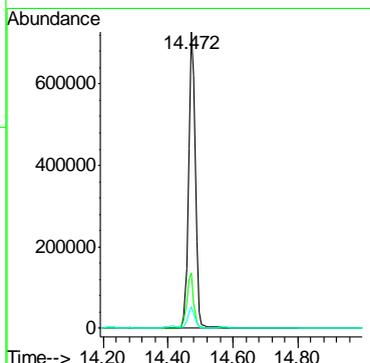
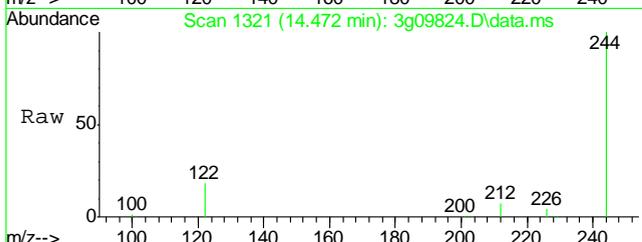
#19  
 Pyrene  
 Concen: 0.0885 ug/mL  
 RT: 14.060 min Scan# 1269  
 Delta R.T. -0.008 min  
 Lab File: 3g09824.D  
 Acq: 22 Jun 12 8:03 pm

Tgt Ion	Resp	Lower	Upper
202	4513		
200	25.0	0.4	40.4
203	21.7	0.0	37.9
201	25.7	0.0	36.8

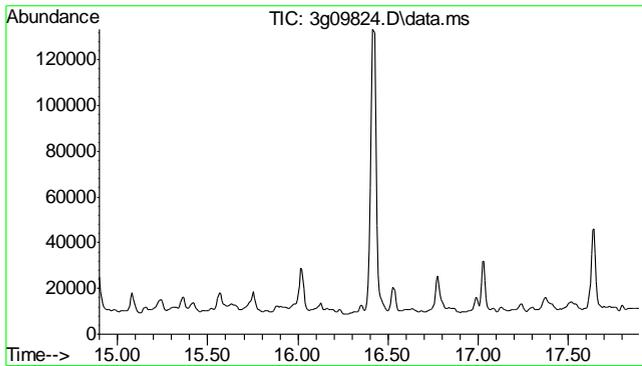


#20  
 Terphenyl-d14  
 Concen: 42.1726 ug/mL  
 RT: 14.472 min Scan# 1321  
 Delta R.T. -0.016 min  
 Lab File: 3g09824.D  
 Acq: 22 Jun 12 8:03 pm

Tgt Ion	Resp	Lower	Upper
244	1071623		
122	18.5	0.0	38.8
212	6.9	0.0	26.9



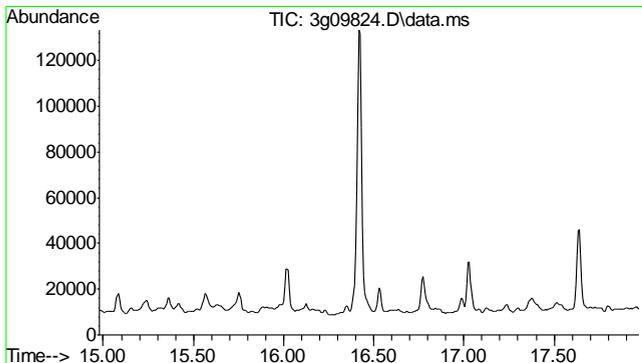
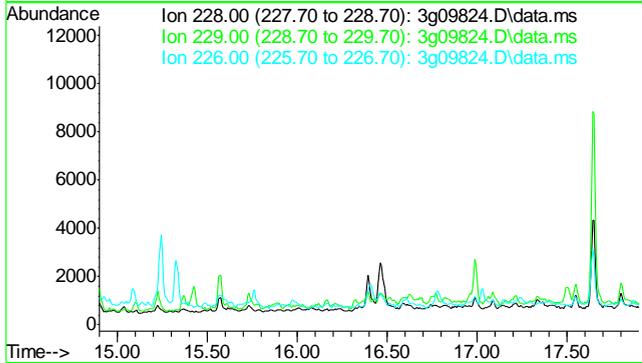
8.1.1  
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#21  
 Benzo(a)anthracene  
 Concen: N.D. ug/mL  
 Expected RT: 16.40 min

Lab File: 3g09824.D  
 Acq: 22 Jun 12 8:03 pm

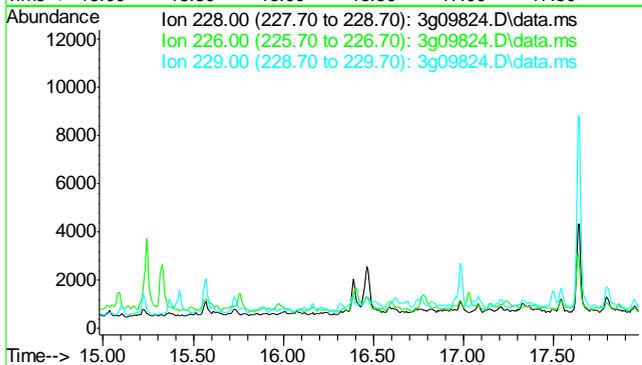
Tgt Ion	Sig	Exp Ratio
228	228	100
229	229	19.4
226	226	26.4

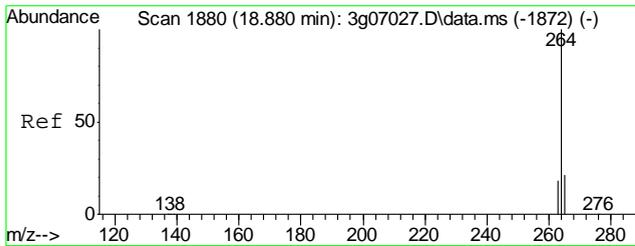


#22  
 Chrysene  
 Concen: N.D. ug/mL  
 Expected RT: 16.48 min

Lab File: 3g09824.D  
 Acq: 22 Jun 12 8:03 pm

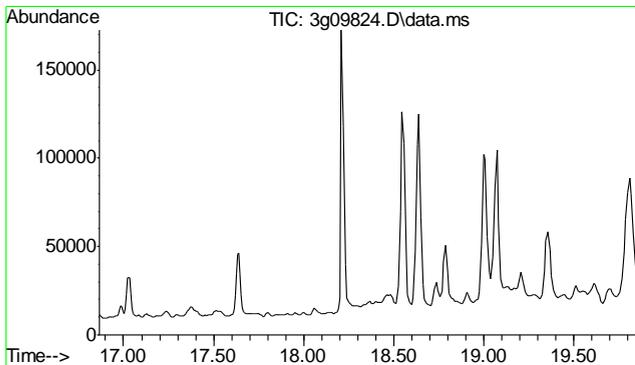
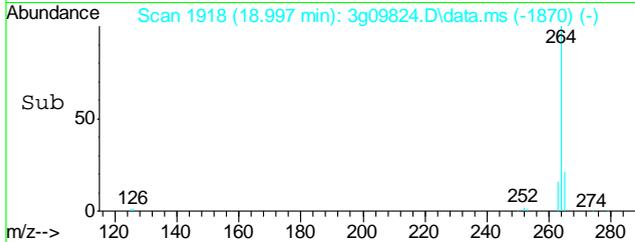
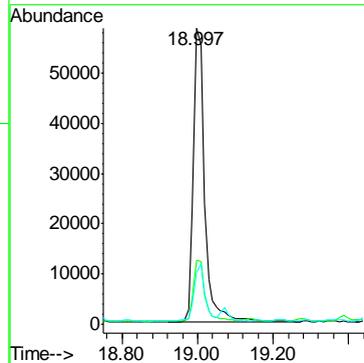
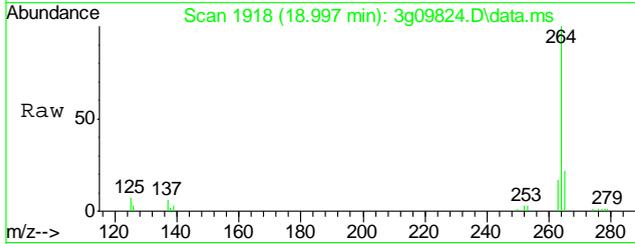
Tgt Ion	Sig	Exp Ratio
228	228	100
226	226	28.9
229	229	19.3





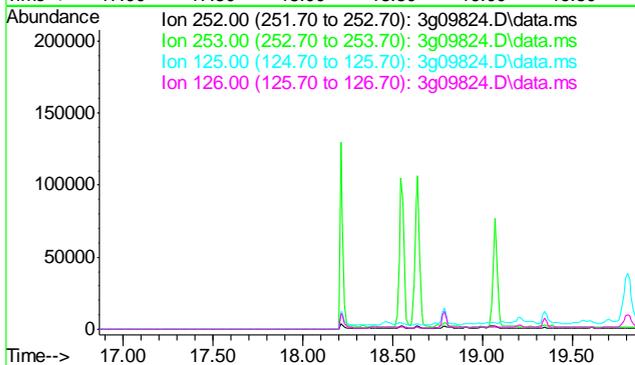
#23  
 Perylene-d12  
 Concen: 4.0000 ug/mL  
 RT: 18.997 min Scan# 1918  
 Delta R.T. 0.000 min  
 Lab File: 3g09824.D  
 Acq: 22 Jun 12 8:03 pm

Tgt Ion	Resp	Lower	Upper
264	120082	100	
265	20.7	1.1	41.1
263	18.9	0.0	38.6

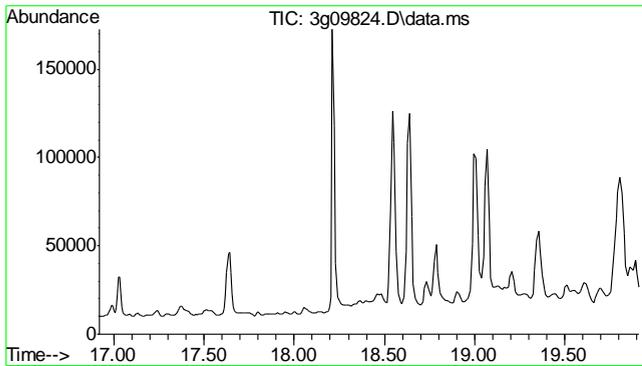


#24  
 Benzo(b)fluoranthene  
 Concen: N.D. ug/mL  
 Expected RT: 18.37 min  
 Lab File: 3g09824.D  
 Acq: 22 Jun 12 8:03 pm

Tgt Ion	Sig	Exp Ratio
252	100	
253	21.3	
125	14.7	
126	20.4	



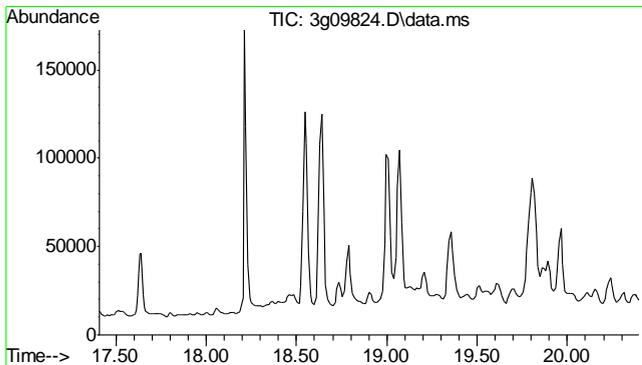
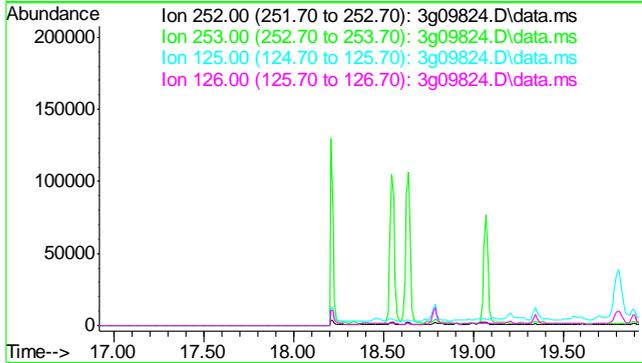
8.1.1  
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#25  
 Benzo(k)fluoranthene  
 Concen: N.D. ug/mL  
 Expected RT: 18.42 min

Lab File: 3g09824.D  
 Acq: 22 Jun 12 8:03 pm

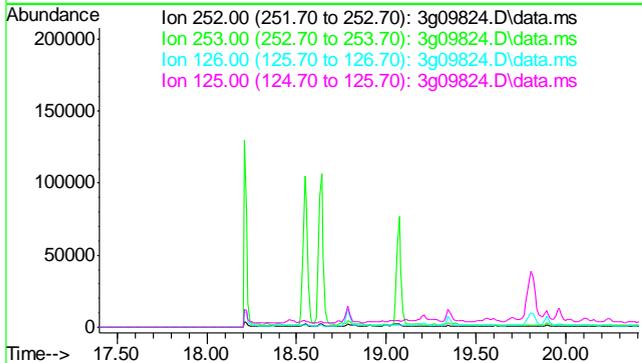
Tgt Ion	Exp Ratio
252	100
253	21.8
125	12.9
126	20.1

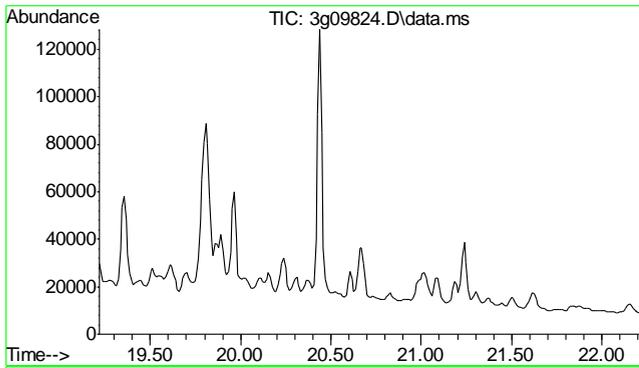


#26  
 Benzo(a)pyrene  
 Concen: N.D. ug/mL  
 Expected RT: 18.90 min

Lab File: 3g09824.D  
 Acq: 22 Jun 12 8:03 pm

Tgt Ion	Exp Ratio
252	100
253	21.6
126	19.7
125	14.8

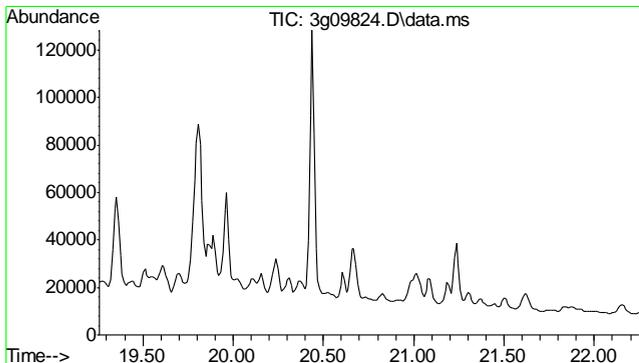
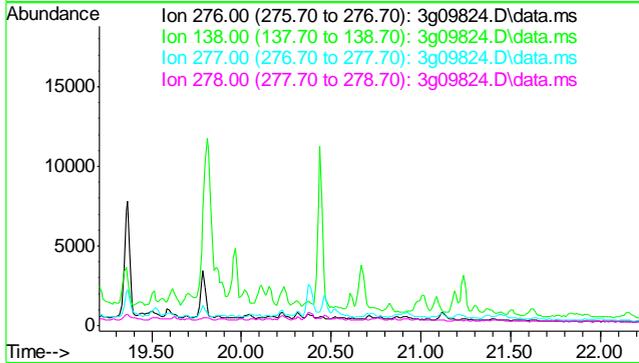




#27  
 Indeno(1,2,3-cd)pyrene  
 Concen: N.D. ug/mL  
 Expected RT: 20.71 min

Lab File: 3g09824.D  
 Acq: 22 Jun 12 8:03 pm

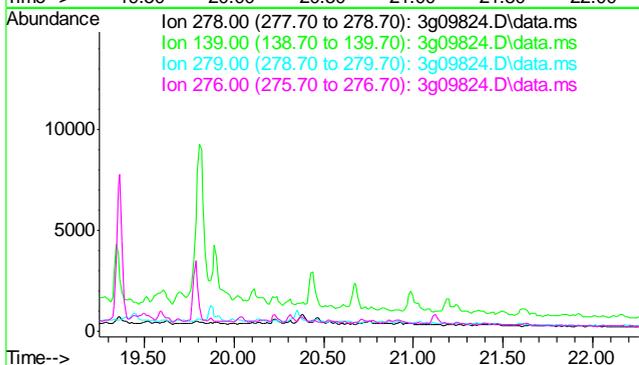
Tgt Ion	Exp Ratio
276	100
138	31.5
277	25.0
278	81.3

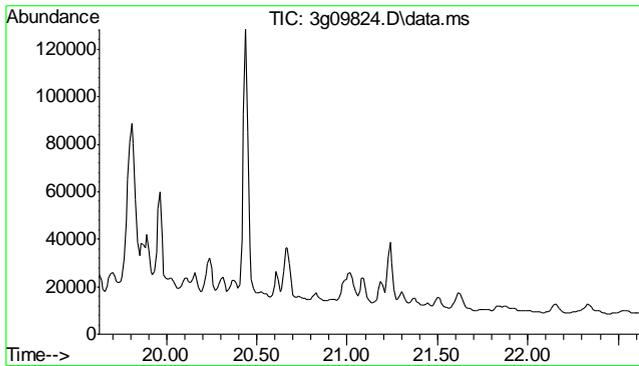


#28  
 Dibenz(a,h)anthracene  
 Concen: N.D. ug/mL  
 Expected RT: 20.75 min

Lab File: 3g09824.D  
 Acq: 22 Jun 12 8:03 pm

Tgt Ion	Exp Ratio
278	100
139	21.3
279	23.4
276	123.0

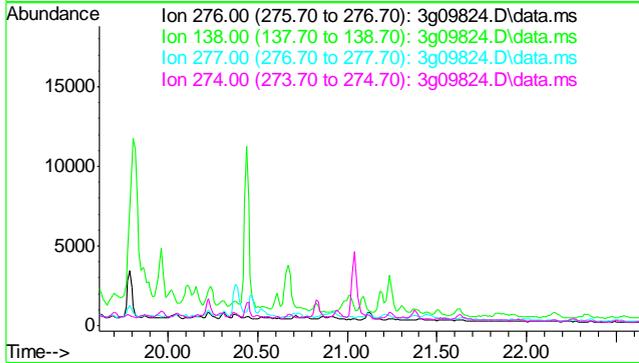




#29  
 Benzo(g,h,i)perylene  
 Concen: N.D. ug/mL  
 Expected RT: 21.12 min

Lab File: 3g09824.D  
 Acq: 22 Jun 12 8:03 pm

Tgt Ion:	276
Sig	Exp Ratio
276	100
138	27.1
277	23.7
274	21.6



8.1.1  
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Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\062212\  
 Data File : 3g09818.D  
 Acq On : 22 Jun 2012 4:19 pm  
 Operator : SARAHM1  
 Sample : OP6113-MB  
 Misc : OP6113,E3G436,30.00,,,1,1  
 ALS Vial : 11 Sample Multiplier: 1

Quant Time: Jun 25 10:17:13 2012  
 Quant Method : C:\msdchem\1\METHODS\SIMPE3G436.M  
 Quant Title : PAHSIM BASE  
 QLast Update : Fri Jun 22 16:07:04 2012  
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Naphthalene-d8	6.429	136	246720	4.0000	ug/mL	0.00
6) Acenaphthene-d10	8.823	164	141681	4.0000	ug/mL	0.00
14) Phenanthrene-d10	11.369	188	221405	4.0000	ug/mL	0.00
18) Chrysene-d12	16.417	240	186931	4.0000	ug/mL	0.00
23) Perylene-d12	18.997	264	102831	4.0000	ug/mL	0.00

System Monitoring Compounds						
2) Nitrobenzene-d5	5.718	82	1341362	37.8914	ug/mL	-0.01
Spiked Amount	50.000	Range 25 - 135	Recovery	=	75.78%	
7) 2-Fluorobiphenyl	7.807	172	1871755	39.7966	ug/mL	-0.01
Spiked Amount	50.000	Range 25 - 135	Recovery	=	79.60%	
20) Terphenyl-d14	14.472	244	1636911	48.3007	ug/mL	-0.02
Spiked Amount	50.000	Range 25 - 135	Recovery	=	96.60%	

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	0.000	128	0	N.D.	d
8) 2-Methylnaphthalene	0.000	142	0	N.D.	d
9) 1-Methylnaphthalene	0.000	142	0	N.D.	d
10) Acenaphthylene	0.000	152	0	N.D.	d
11) Acenaphthene	0.000	154	0	N.D.	d
12) Fluorene	0.000	166	0	N.D.	d
13) Diphenylamine	0.000	169	0	N.D.	d
15) Phenanthrene	0.000	178	0	N.D.	d
16) Anthracene	0.000	178	0	N.D.	d
17) Fluoranthene	0.000	202	0	N.D.	d
19) Pyrene	0.000	202	0	N.D.	d
21) Benzo(a)anthracene	0.000	228	0	N.D.	d
22) Chrysene	0.000	228	0	N.D.	d
24) Benzo(b)fluoranthene	0.000	252	0	N.D.	d
25) Benzo(k)fluoranthene	0.000	252	0	N.D.	d
26) Benzo(a)pyrene	0.000	252	0	N.D.	d
27) Indeno(1,2,3-cd)pyrene	0.000	276	0	N.D.	d
28) Dibenzo(a,h)anthracene	0.000	278	0	N.D.	d
29) Benzo(g,h,i)perylene	0.000	276	0	N.D.	d

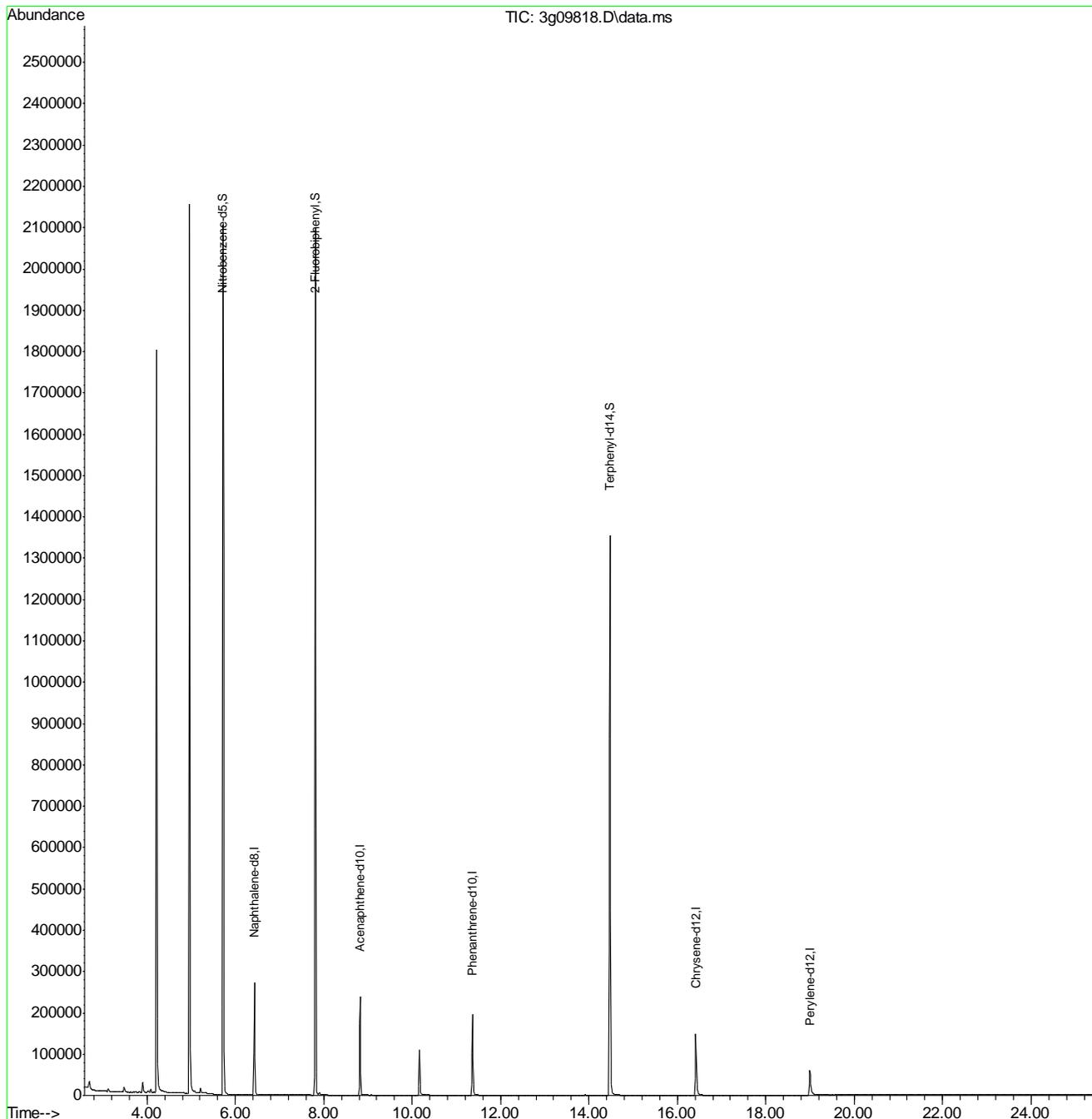
(#) = qualifier out of range (m) = manual integration (+) = signals summed

8.2.1  
 8

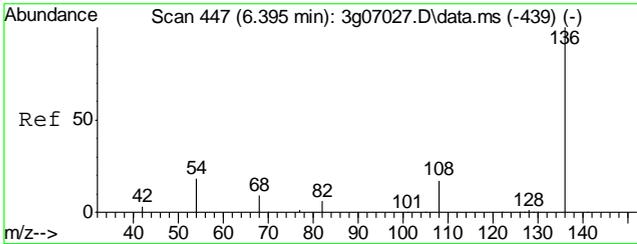
Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\062212\  
 Data File : 3g09818.D  
 Acq On : 22 Jun 2012 4:19 pm  
 Operator : SARAHM1  
 Sample : OP6113-MB  
 Misc : OP6113,E3G436,30.00,,,1,1  
 ALS Vial : 11 Sample Multiplier: 1

Quant Time: Jun 25 10:17:13 2012  
 Quant Method : C:\msdchem\1\METHODS\SIMPE3G436.M  
 Quant Title : PAHSIM BASE  
 QLast Update : Fri Jun 22 16:07:04 2012  
 Response via : Initial Calibration

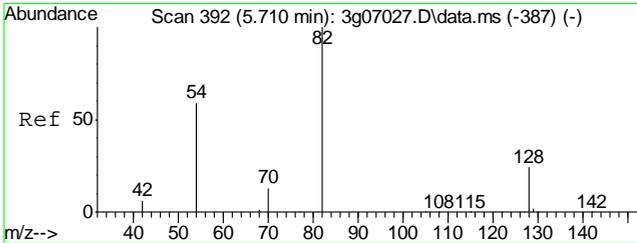
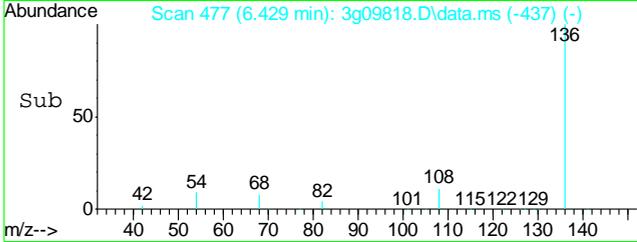
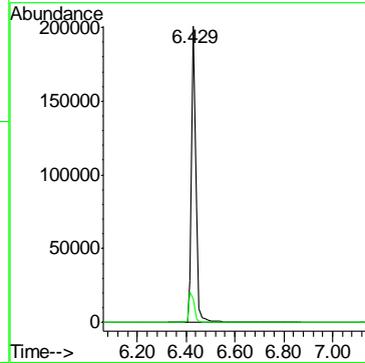
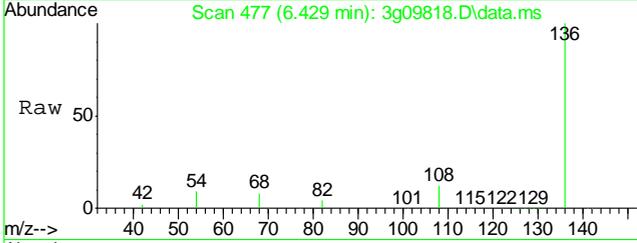


8.2.1  
8



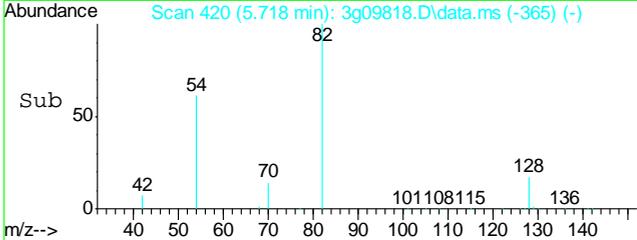
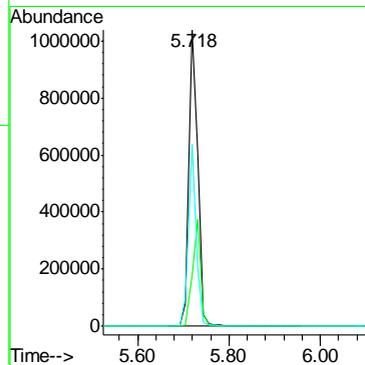
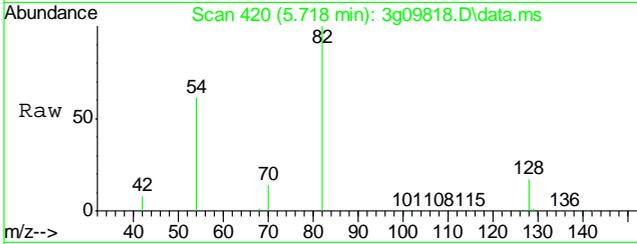
#1  
 Naphthalene-d8  
 Concen: 4.0000 ug/mL  
 RT: 6.429 min Scan# 477  
 Delta R.T. 0.000 min  
 Lab File: 3g09818.D  
 Acq: 22 Jun 12 4:19 pm

Tgt Ion	Resp	Lower	Upper
136	246720	100	
68	12.4	0.0	32.0

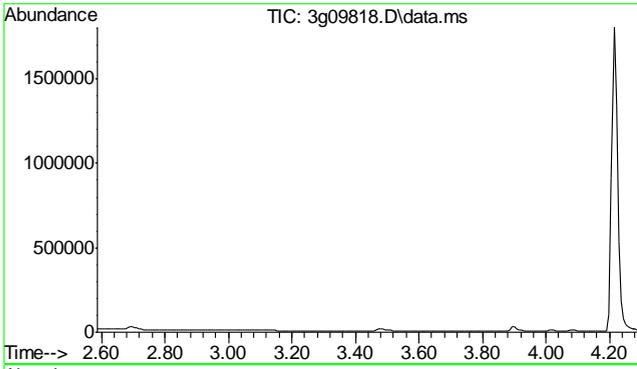


#2  
 Nitrobenzene-d5  
 Concen: 37.8914 ug/mL  
 RT: 5.718 min Scan# 420  
 Delta R.T. -0.012 min  
 Lab File: 3g09818.D  
 Acq: 22 Jun 12 4:19 pm

Tgt Ion	Resp	Lower	Upper
82	1341362	100	
128	33.8	15.2	55.2
54	54.1	34.0	74.0



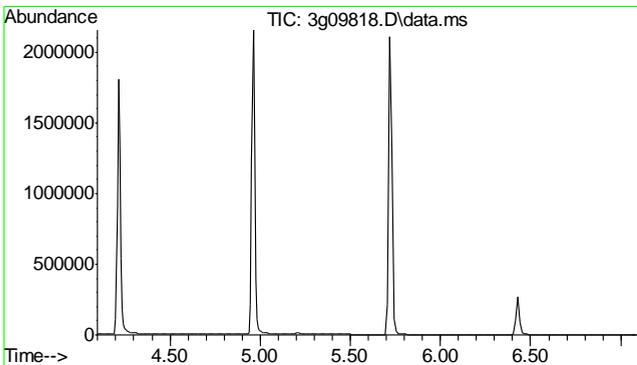
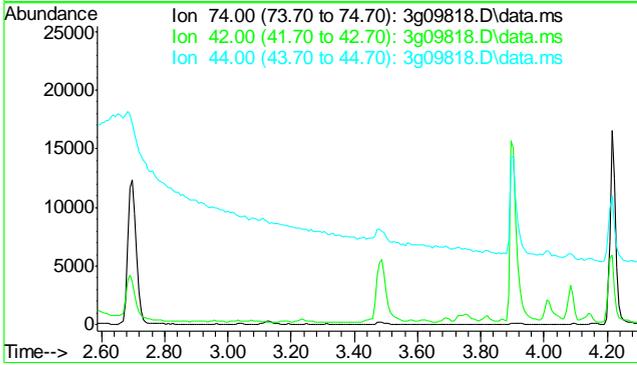
8.2.1  
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#3  
 N-Nitrosodimethylamine  
 Concen: N.D. ug/mL  
 Expected RT: 2.79 min

Lab File: 3g09818.D  
 Acq: 22 Jun 12 4:19 pm

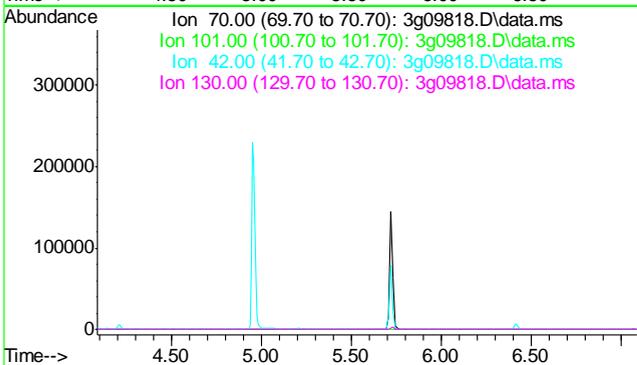
Tgt Ion:	74
Sig	Exp Ratio
74	100
42	51.8
44	10.0

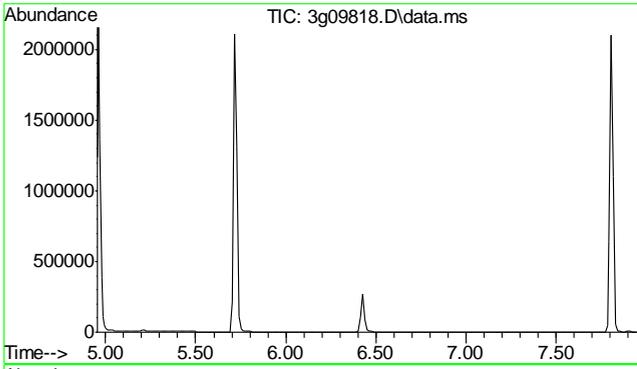


#4  
 N-Nitrosodi-propylamine  
 Concen: N.D. ug/mL  
 Expected RT: 5.59 min

Lab File: 3g09818.D  
 Acq: 22 Jun 12 4:19 pm

Tgt Ion:	70
Sig	Exp Ratio
70	100
101	10.3
42	59.5
130	20.1



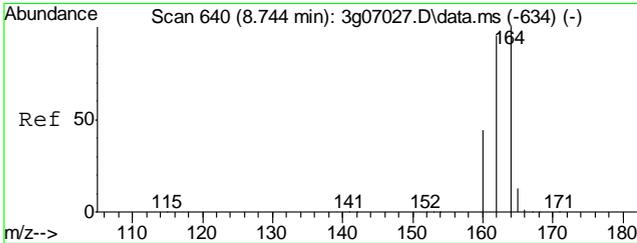
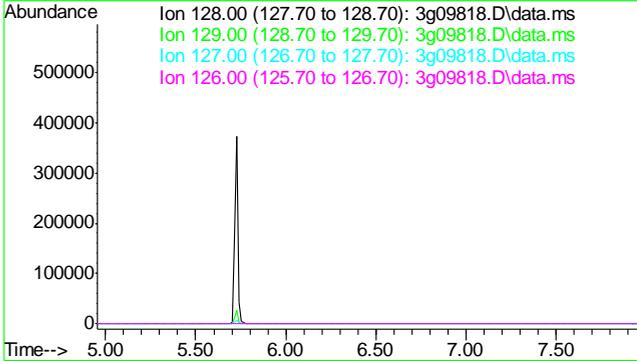


#5  
 Naphthalene  
 Concen: N.D. ug/mL  
 Expected RT: 6.45 min

Lab File: 3g09818.D  
 Acq: 22 Jun 12 4:19 pm

Tgt Ion: 128

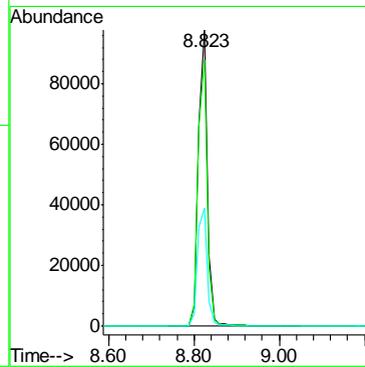
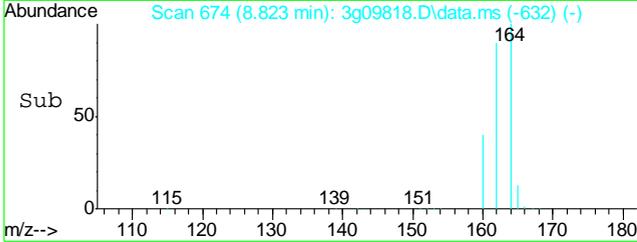
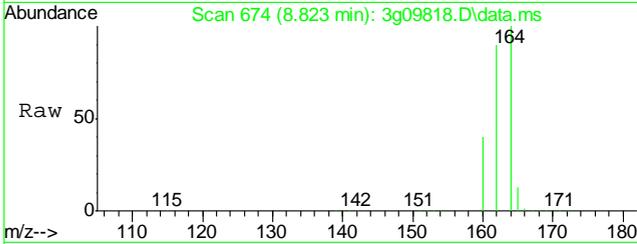
Sig	Exp Ratio
128	100
129	10.9
127	12.4
126	7.7

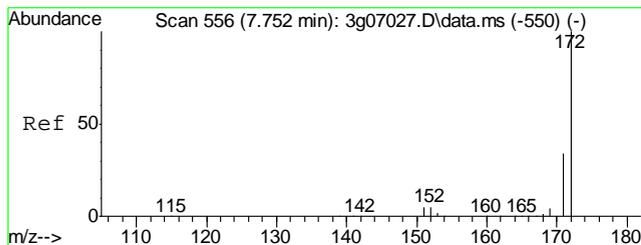


#6  
 Acenaphthene-d10  
 Concen: 4.0000 ug/mL  
 RT: 8.823 min Scan# 674  
 Delta R.T. 0.000 min  
 Lab File: 3g09818.D  
 Acq: 22 Jun 12 4:19 pm

Tgt Ion: 164 Resp: 141681

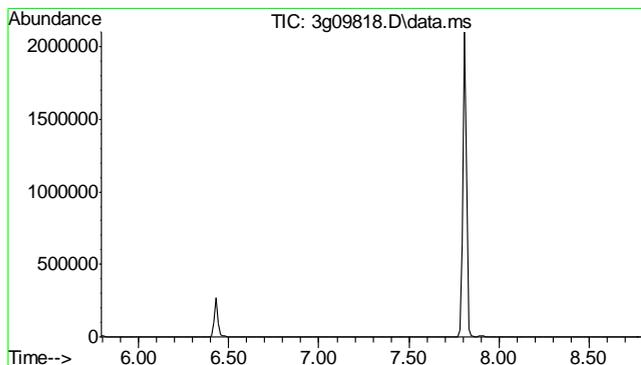
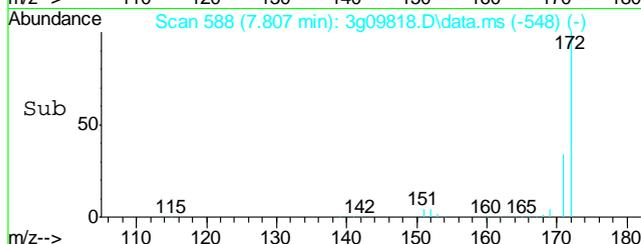
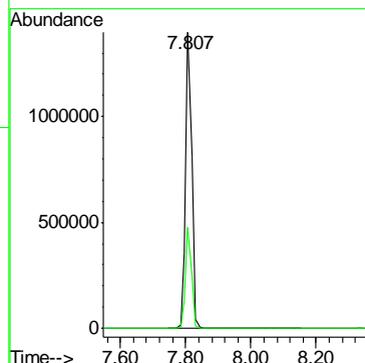
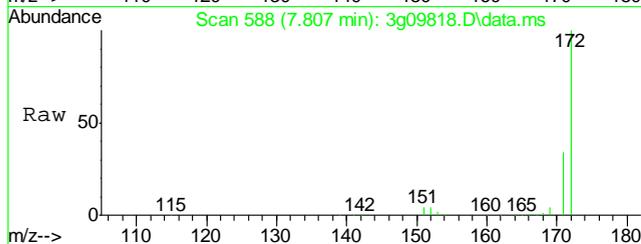
Ion	Ratio	Lower	Upper
164	100		
162	93.3	72.5	112.5
160	43.2	20.9	60.9





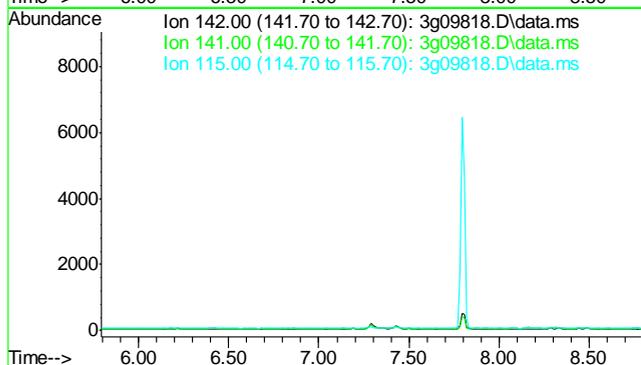
#7  
 2-Fluorobiphenyl  
 Concen: 39.7966 ug/mL  
 RT: 7.807 min Scan# 588  
 Delta R.T. -0.012 min  
 Lab File: 3g09818.D  
 Acq: 22 Jun 12 4:19 pm

Tgt Ion: 172 Resp: 1871755  
 Ion Ratio Lower Upper  
 172 100  
 171 33.2 13.2 53.2

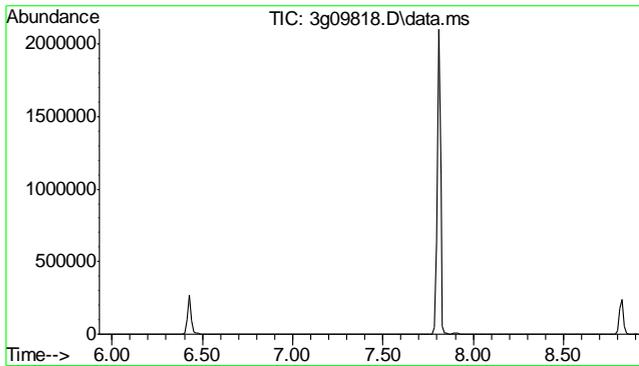


#8  
 2-Methylnaphthalene  
 Concen: N.D. ug/mL  
 Expected RT: 7.29 min  
 Lab File: 3g09818.D  
 Acq: 22 Jun 12 4:19 pm

Tgt Ion: 142  
 Sig Exp Ratio  
 142 100  
 141 83.1  
 115 34.8

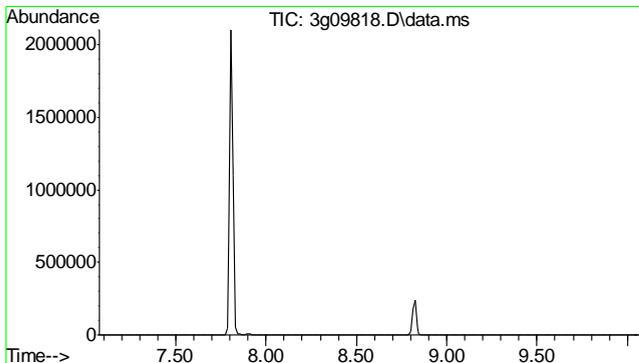
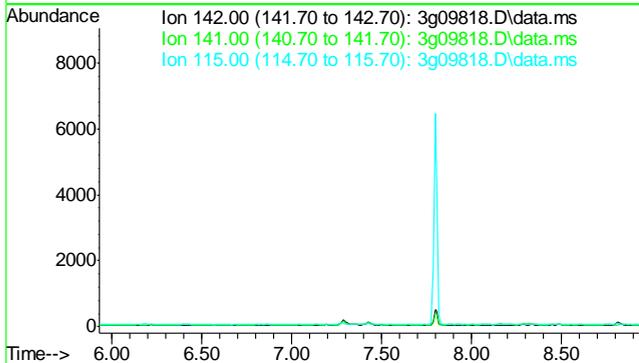


8.2.1  
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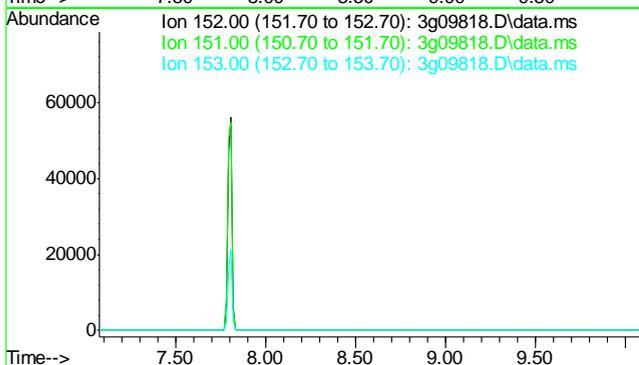
#9  
 1-Methylnaphthalene  
 Concen: N.D. ug/mL  
 Expected RT: 7.43 min  
 Lab File: 3g09818.D  
 Acq: 22 Jun 12 4:19 pm

Tgt Ion	Sig	Exp Ratio
142	142	100
141	141	86.4
115	115	36.1

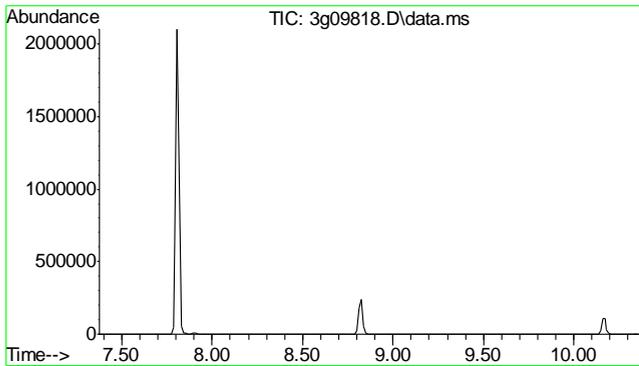


#10  
 Acenaphthylene  
 Concen: N.D. ug/mL  
 Expected RT: 8.58 min  
 Lab File: 3g09818.D  
 Acq: 22 Jun 12 4:19 pm

Tgt Ion	Sig	Exp Ratio
152	152	100
151	151	19.1
153	153	14.1

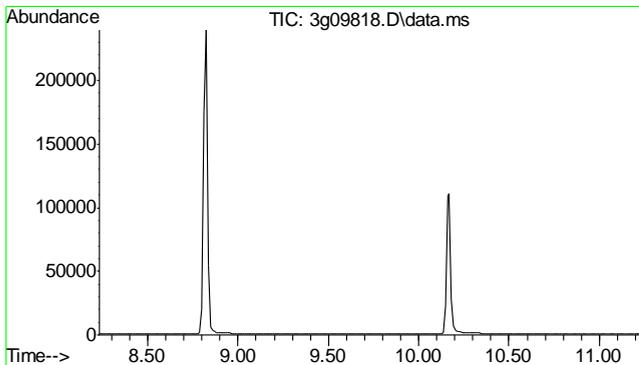
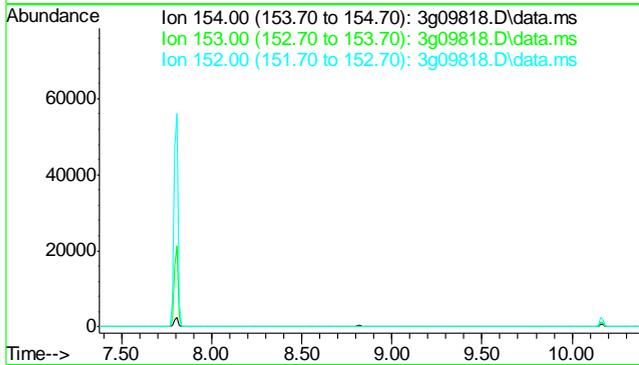


8.2.1  
 8



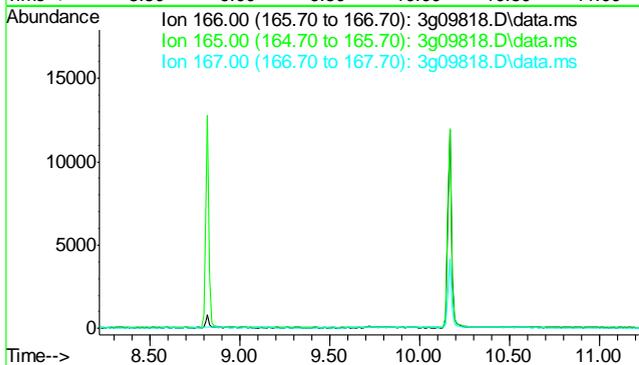
#11  
 Acenaphthene  
 Concen: N.D. ug/mL  
 Expected RT: 8.87 min  
 Lab File: 3g09818.D  
 Acq: 22 Jun 12 4:19 pm

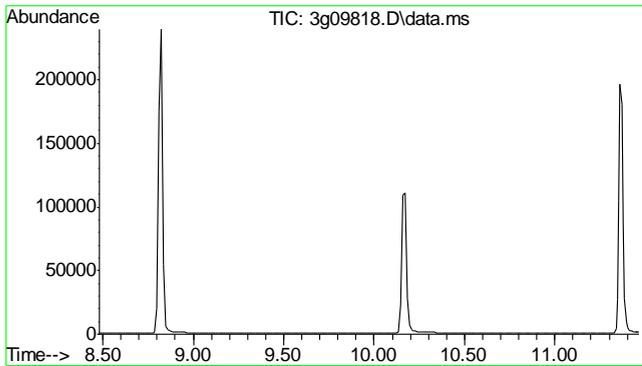
Tgt Ion	Sig	Exp Ratio
154	154	100
153	153	104.2
152	152	45.3



#12  
 Fluorene  
 Concen: N.D. ug/mL  
 Expected RT: 9.72 min  
 Lab File: 3g09818.D  
 Acq: 22 Jun 12 4:19 pm

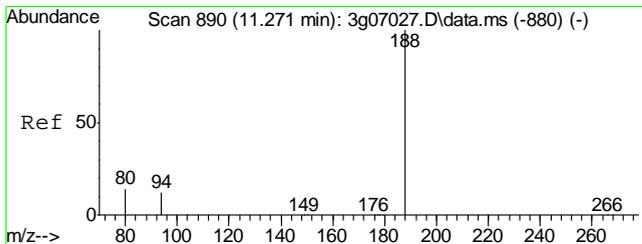
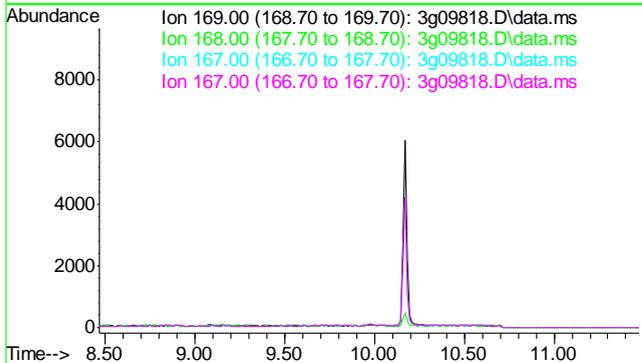
Tgt Ion	Sig	Exp Ratio
166	166	100
165	165	90.8
167	167	13.3





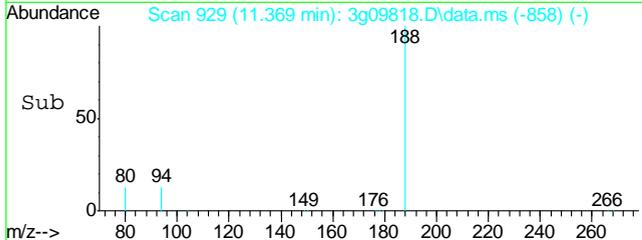
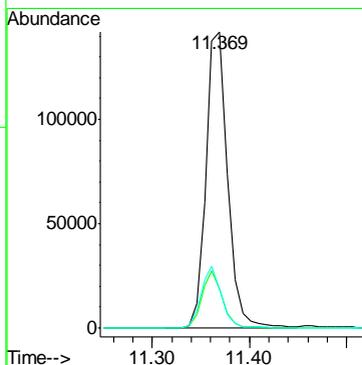
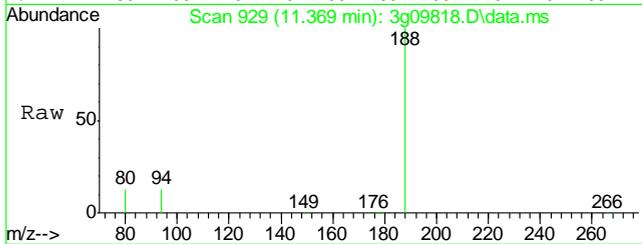
#13  
 Diphenylamine  
 Concen: N.D. ug/mL  
 Expected RT: 9.97 min  
 Lab File: 3g09818.D  
 Acq: 22 Jun 12 4:19 pm

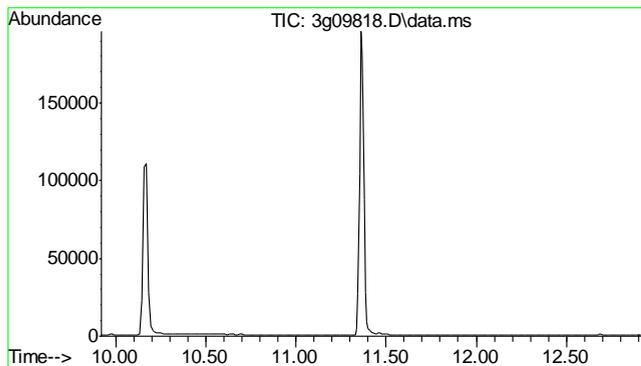
Tgt Ion	Exp Ratio
169	100
168	62.1
167	33.2
167	33.2



#14  
 Phenanthrene-d10  
 Concen: 4.0000 ug/mL  
 RT: 11.369 min Scan# 929  
 Delta R.T. 0.000 min  
 Lab File: 3g09818.D  
 Acq: 22 Jun 12 4:19 pm

Tgt Ion	Resp	Ion Ratio	Lower	Upper
188	221405	100		
94		18.2	0.0	37.2
80		20.0	0.0	38.3

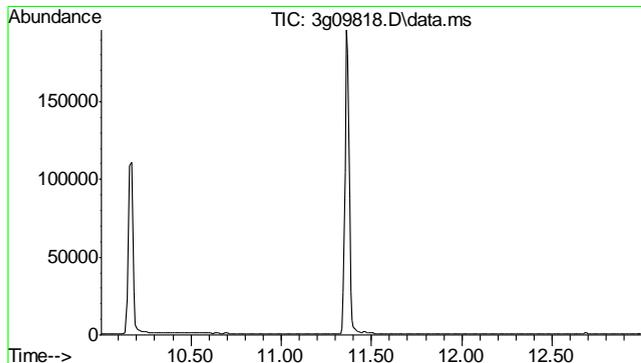
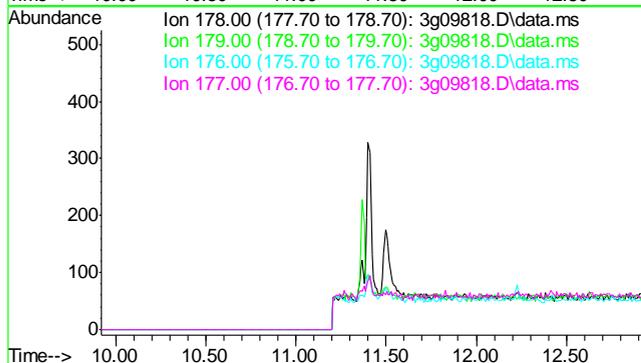




#15  
 Phenanthrene  
 Concen: N.D. ug/mL  
 Expected RT: 11.42 min

Lab File: 3g09818.D  
 Acq: 22 Jun 12 4:19 pm

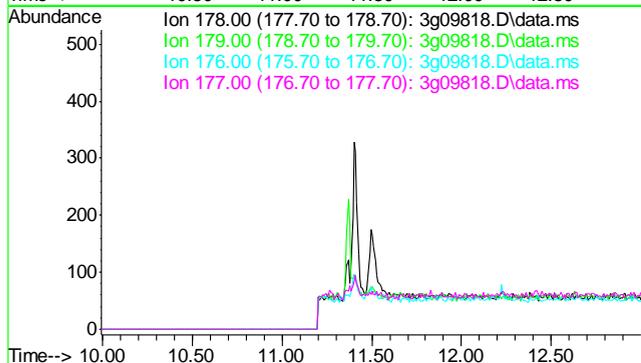
Tgt Ion:	178
Sig	Exp Ratio
178	100
179	15.1
176	18.6
177	10.4

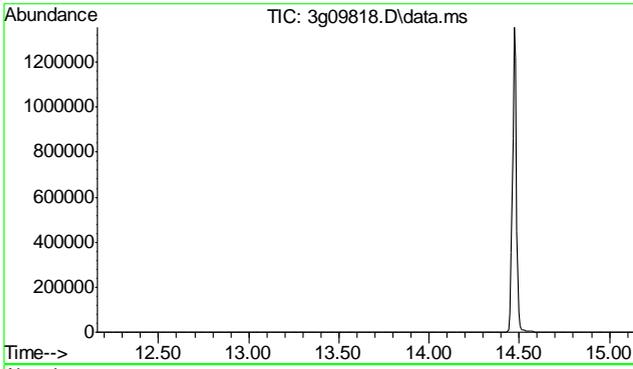


#16  
 Anthracene  
 Concen: N.D. ug/mL  
 Expected RT: 11.50 min

Lab File: 3g09818.D  
 Acq: 22 Jun 12 4:19 pm

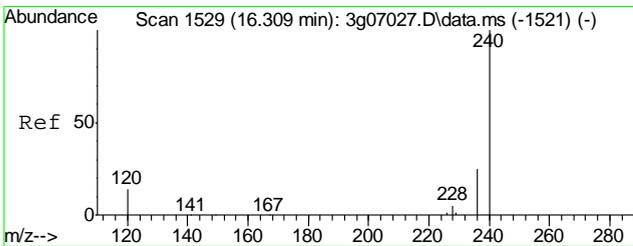
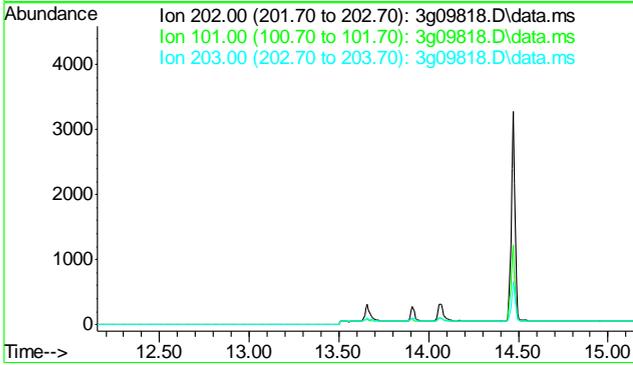
Tgt Ion:	178
Sig	Exp Ratio
178	100
179	15.0
176	17.9
177	8.9





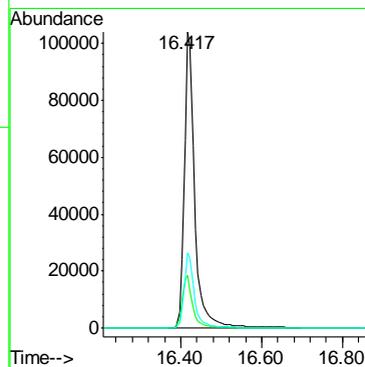
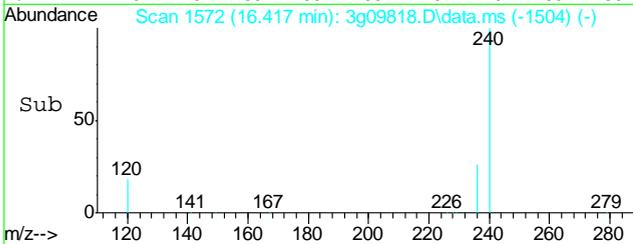
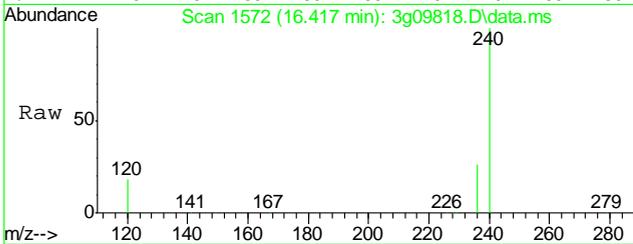
#17  
 Fluoranthene  
 Concen: N.D. ug/mL  
 Expected RT: 13.66 min  
  
 Lab File: 3g09818.D  
 Acq: 22 Jun 12 4:19 pm  
  
 Tgt Ion: 202  

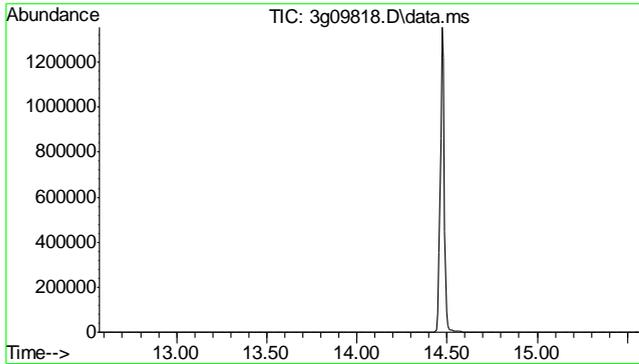
Sig	Exp Ratio
202	100
101	17.6
203	17.1



#18  
 Chrysene-d12  
 Concen: 4.0000 ug/mL  
 RT: 16.417 min Scan# 1572  
 Delta R.T. -0.006 min  
 Lab File: 3g09818.D  
 Acq: 22 Jun 12 4:19 pm

Tgt Ion: 240	Resp: 186931
Ion Ratio	Lower Upper
240	100
120	17.4 0.0 37.2
236	25.3 5.2 45.2

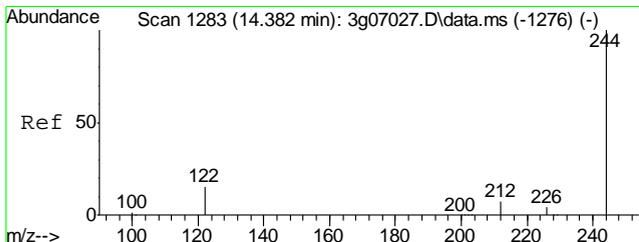
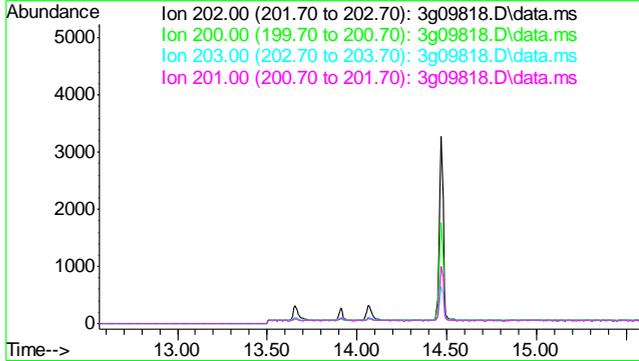




#19  
 Pyrene  
 Concen: N.D. ug/mL  
 Expected RT: 14.07 min

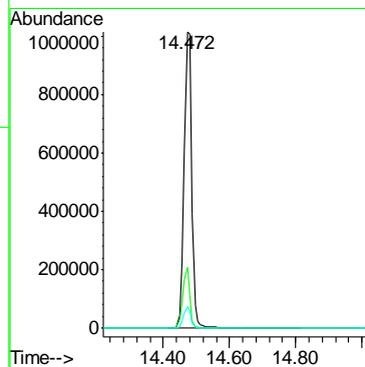
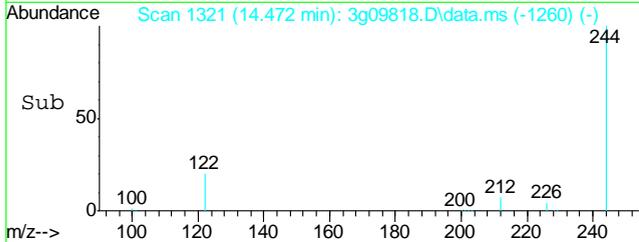
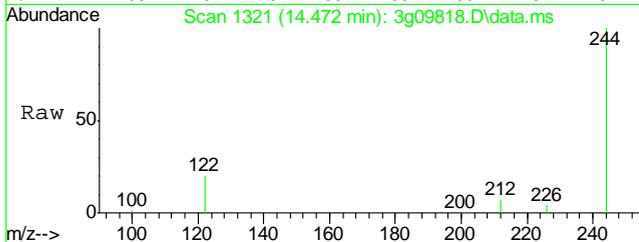
Lab File: 3g09818.D  
 Acq: 22 Jun 12 4:19 pm

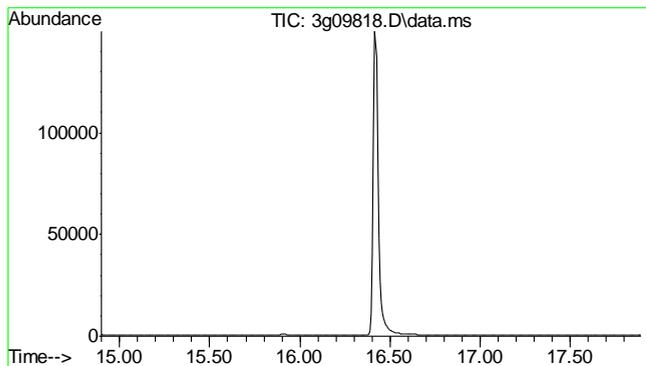
Tgt Ion	Exp Ratio
202	100
200	20.4
203	17.9
201	16.8



#20  
 Terphenyl-d14  
 Concen: 48.3007 ug/mL  
 RT: 14.472 min Scan# 1321  
 Delta R.T. -0.016 min  
 Lab File: 3g09818.D  
 Acq: 22 Jun 12 4:19 pm

Tgt Ion	Resp	Ion Ratio	Lower	Upper
244	1636911	100		
122		18.6	0.0	38.8
212		6.8	0.0	26.9

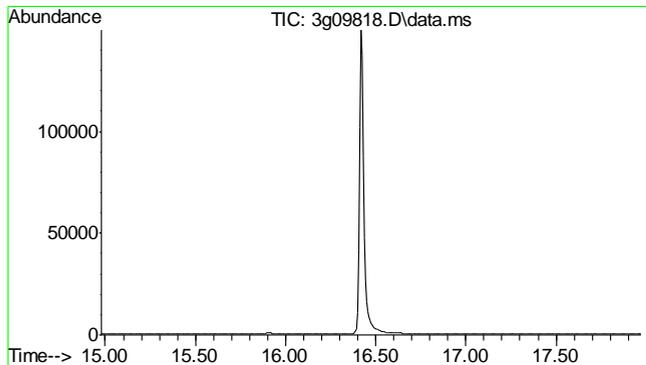
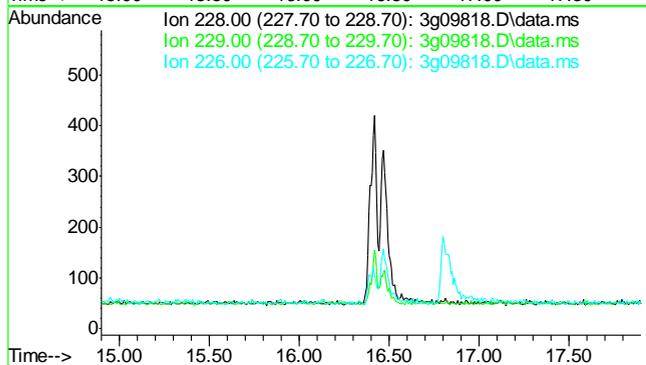




#21  
 Benzo(a)anthracene  
 Concen: N.D. ug/mL  
 Expected RT: 16.40 min

Lab File: 3g09818.D  
 Acq: 22 Jun 12 4:19 pm

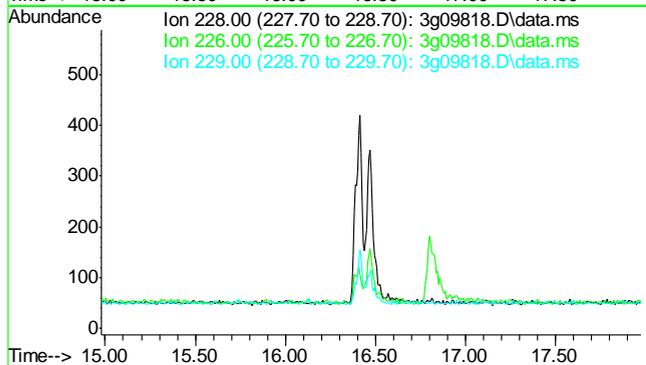
Tgt Ion	Exp Ratio
228	100
229	19.4
226	26.4



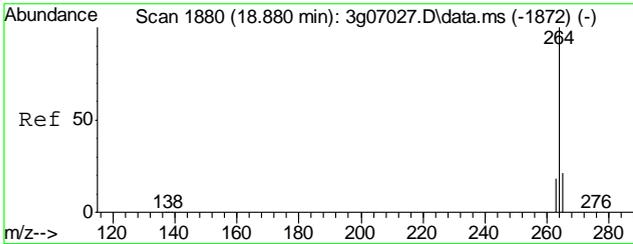
#22  
 Chrysene  
 Concen: N.D. ug/mL  
 Expected RT: 16.48 min

Lab File: 3g09818.D  
 Acq: 22 Jun 12 4:19 pm

Tgt Ion	Exp Ratio
228	100
226	28.9
229	19.3

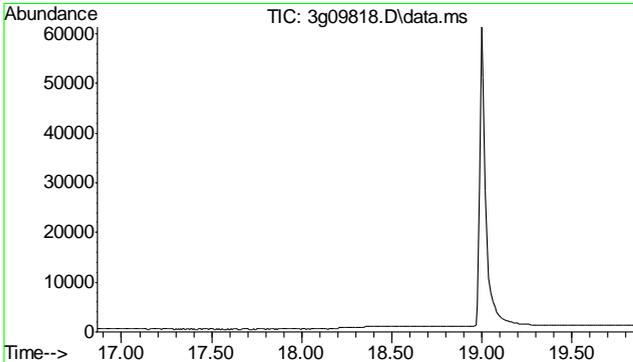
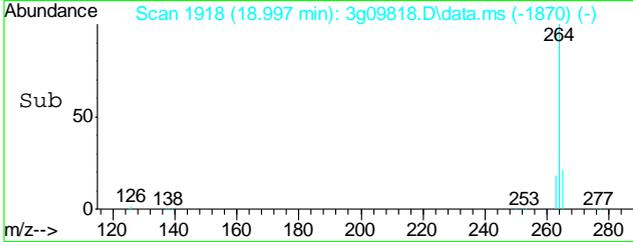
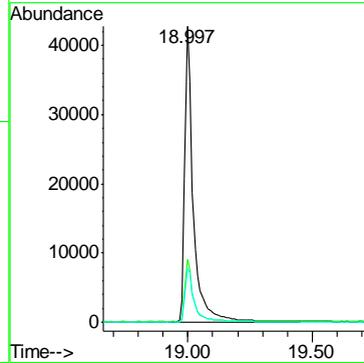
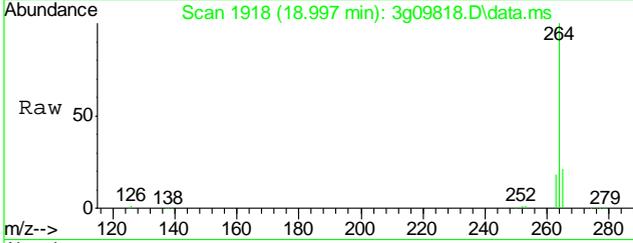


8.2.1  
 8



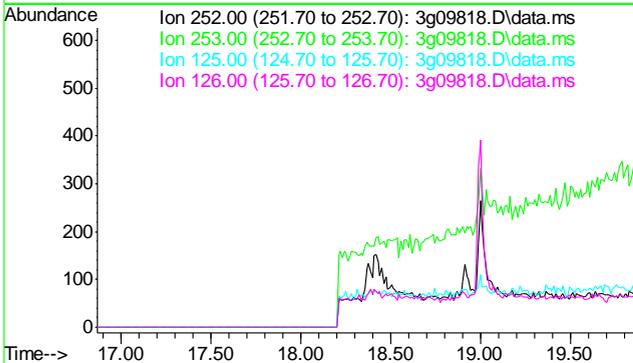
#23  
 Perylene-d12  
 Concen: 4.0000 ug/mL  
 RT: 18.997 min Scan# 1918  
 Delta R.T. 0.000 min  
 Lab File: 3g09818.D  
 Acq: 22 Jun 12 4:19 pm

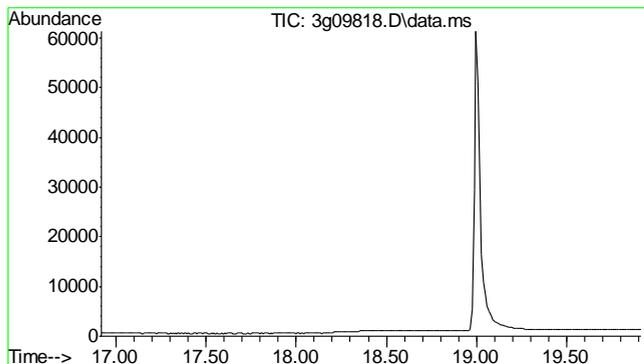
Tgt Ion	Resp	Lower	Upper
264	102831		
265	20.6	1.1	41.1
263	19.4	0.0	38.6



#24  
 Benzo(b)fluoranthene  
 Concen: N.D. ug/mL  
 Expected RT: 18.37 min  
 Lab File: 3g09818.D  
 Acq: 22 Jun 12 4:19 pm

Tgt Ion	Exp Ratio
252	100
253	21.3
125	14.7
126	20.4

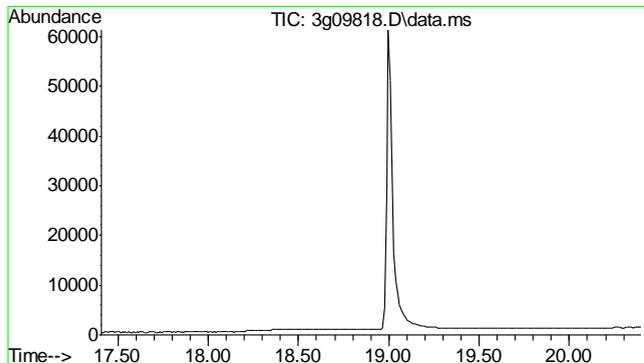
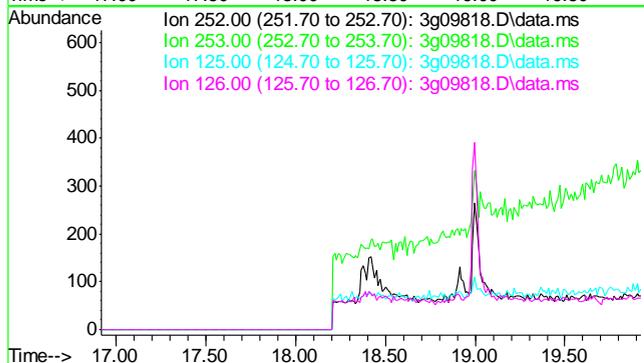




#25  
 Benzo(k)fluoranthene  
 Concen: N.D. ug/mL  
 Expected RT: 18.42 min

Lab File: 3g09818.D  
 Acq: 22 Jun 12 4:19 pm

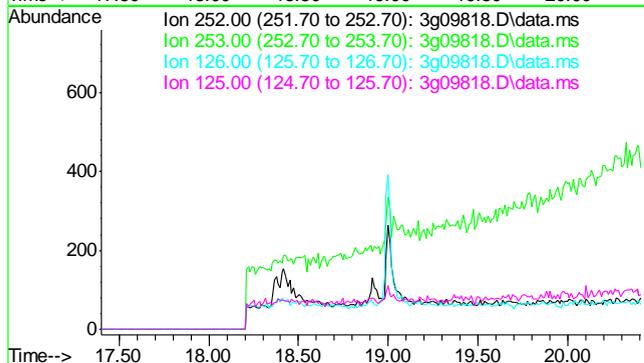
Tgt Ion	Exp Ratio
252	100
253	21.8
125	12.9
126	20.1

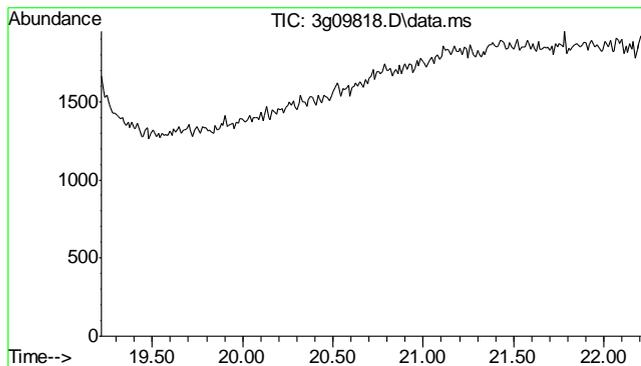


#26  
 Benzo(a)pyrene  
 Concen: N.D. ug/mL  
 Expected RT: 18.90 min

Lab File: 3g09818.D  
 Acq: 22 Jun 12 4:19 pm

Tgt Ion	Exp Ratio
252	100
253	21.6
126	19.7
125	14.8

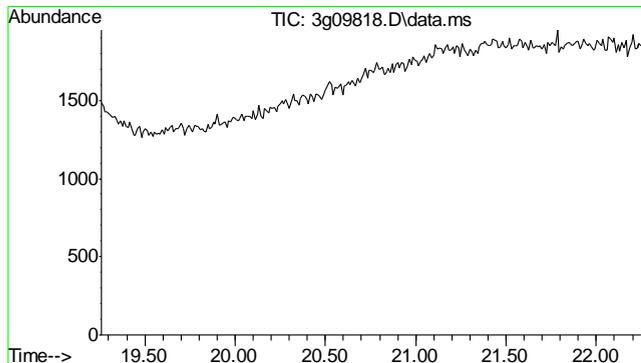
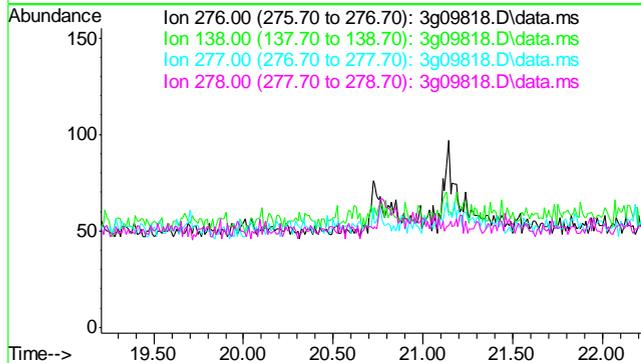




#27  
 Indeno(1,2,3-cd)pyrene  
 Concen: N.D. ug/mL  
 Expected RT: 20.71 min

Lab File: 3g09818.D  
 Acq: 22 Jun 12 4:19 pm

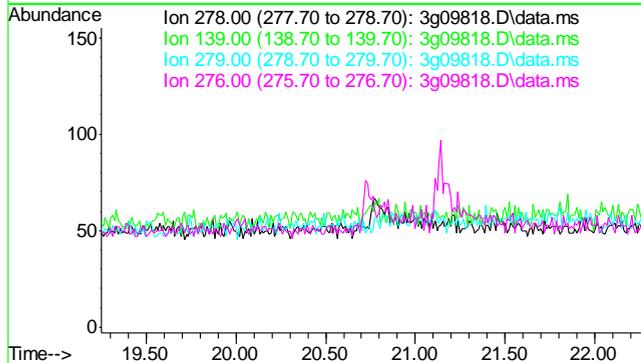
Tgt Ion:	276
Sig	Exp Ratio
276	100
138	31.5
277	25.0
278	81.3



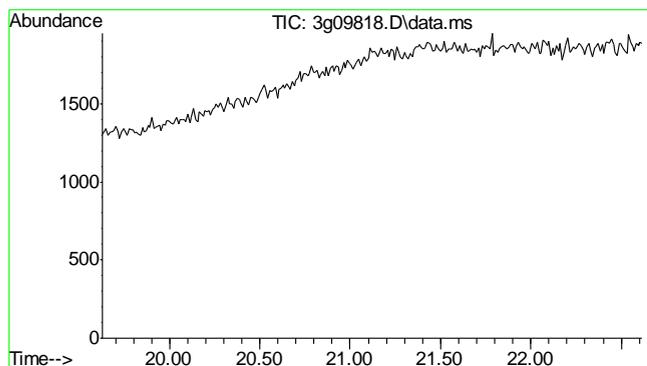
#28  
 Dibenz(a,h)anthracene  
 Concen: N.D. ug/mL  
 Expected RT: 20.75 min

Lab File: 3g09818.D  
 Acq: 22 Jun 12 4:19 pm

Tgt Ion:	278
Sig	Exp Ratio
278	100
139	21.3
279	23.4
276	123.0



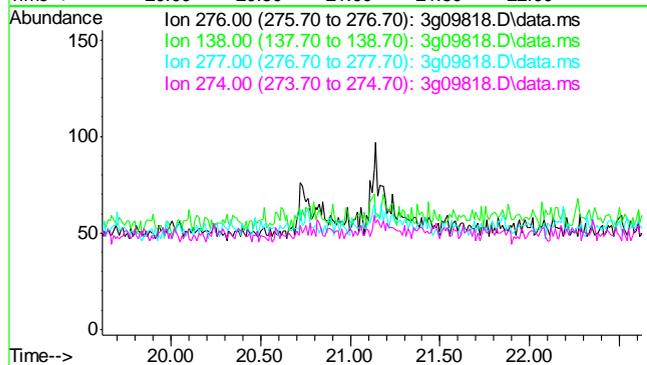
8.2.1  
 8



#29  
 Benzo(g,h,i)perylene  
 Concen: N.D. ug/mL  
 Expected RT: 21.12 min

Lab File: 3g09818.D  
 Acq: 22 Jun 12 4:19 pm

Tgt Ion:	276
Sig	Exp Ratio
276	100
138	27.1
277	23.7
274	21.6



8.2.1  
 8

## 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:** D35710  
**Account:** XTOKRWR XTO Energy  
**Project:** FRU 297-8B

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
GGB910-MB	GB16414.D	1	06/21/12	SK	n/a	n/a	GGB910

The QC reported here applies to the following samples:

Method: SW846 8015B

D35710-1

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

CAS No.	Surrogate Recoveries	Limits
120-82-1	1,2,4-Trichlorobenzene	100% 60-140%

# Blank Spike Summary

**Job Number:** D35710  
**Account:** XTOKRWR XTO Energy  
**Project:** FRU 297-8B

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
GGB910-BS	GB16415.D	1	06/21/12	SK	n/a	n/a	GGB910

The QC reported here applies to the following samples:

Method: SW846 8015B

D35710-1

CAS No.	Compound	Spike mg/kg	BSP mg/kg	BSP %	Limits
	TPH-GRO (C6-C10)	110	111	101	70-130

CAS No.	Surrogate Recoveries	BSP	Limits
120-82-1	1,2,4-Trichlorobenzene	113%	60-140%

9.2.1  
9

\* = Outside of Control Limits.

# Matrix Spike/Matrix Spike Duplicate Summary

**Job Number:** D35710  
**Account:** XTOKRWR XTO Energy  
**Project:** FRU 297-8B

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
D35708-1MS	GB16417.D	1	06/21/12	SK	n/a	n/a	GGB910
D35708-1MSD	GB16418.D	1	06/21/12	SK	n/a	n/a	GGB910
D35708-1	GB16416.D	1	06/21/12	SK	n/a	n/a	GGB910

The QC reported here applies to the following samples:

Method: SW846 8015B

D35710-1

CAS No.	Compound	D35708-1 mg/kg	Spike Q	mg/kg	MS mg/kg	MS %	MSD mg/kg	MSD %	RPD	Limits Rec/RPD
	TPH-GRO (C6-C10)	ND	121	125	103	124	102	1	70-130/30	

CAS No.	Surrogate Recoveries	MS	MSD	D35708-1	Limits
120-82-1	1,2,4-Trichlorobenzene	106%	105%	98%	60-140%

9.3.1  
9

\* = Outside of Control Limits.

GC Volatiles

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Raw Data

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Manual Integrations  
**APPROVED**  
 (compounds with "m" flag)  
**Judy Melson**  
**06/25/12 09:41**

Quantitation Report (QT Reviewed)

Signal #1 : Y:\1\DATA\062112\GB16420.D\FID1A.CH Vial: 9  
 Signal #2 : Y:\1\DATA\062112\GB16420.D\FID2B.CH  
 Acq On : 21 Jun 2012 7:05 pm Operator: StephK  
 Sample : D35710-1, 50X Inst : GC/MS Ins  
 Misc : GC2928,GGB910,5.044,,100,5,1 Multiplr: 1.00  
 IntFile Signal #1: TVH1.E IntFile Signal #2: FB2.E  
 Quant Time: Jun 22 08:09:10 2012 Quant Results File: TB868GB868SOIL.RES

Quant Method : C:\MSDCHEM\1...\TB868GB868SOIL.M (Chemstation Integrator)  
 Title : 8015B/8021B TVH/BTEX  
 Last Update : Thu Jun 21 15:17:12 2012  
 Response via : Initial Calibration  
 DataAcq Meth : TVB4.M

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

Compound	R.T.	Response	Conc	Units
System Monitoring Compounds				
2) S 1,2,4-Trichlorobenzene	14.36	2982648	95.189 %	m
10) S 1,2,4-Trichlorobenzene (P)	14.36	17634173	108.500 %	
Target Compounds				
1) H TVH-Gasoline	7.23	5980887	<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	7.63	244379	0.617	ug/L
7) T Ethylbenzene	10.27	95381	0.282	ug/L
8) T m,p-Xylene	10.45	407218	0.742	ug/L
9) T o-Xylene	0.00	0	N.D.	ug/L d
11) T Naphthalene	14.54	1354212	6.863	ug/L

10.1.1  
**10**

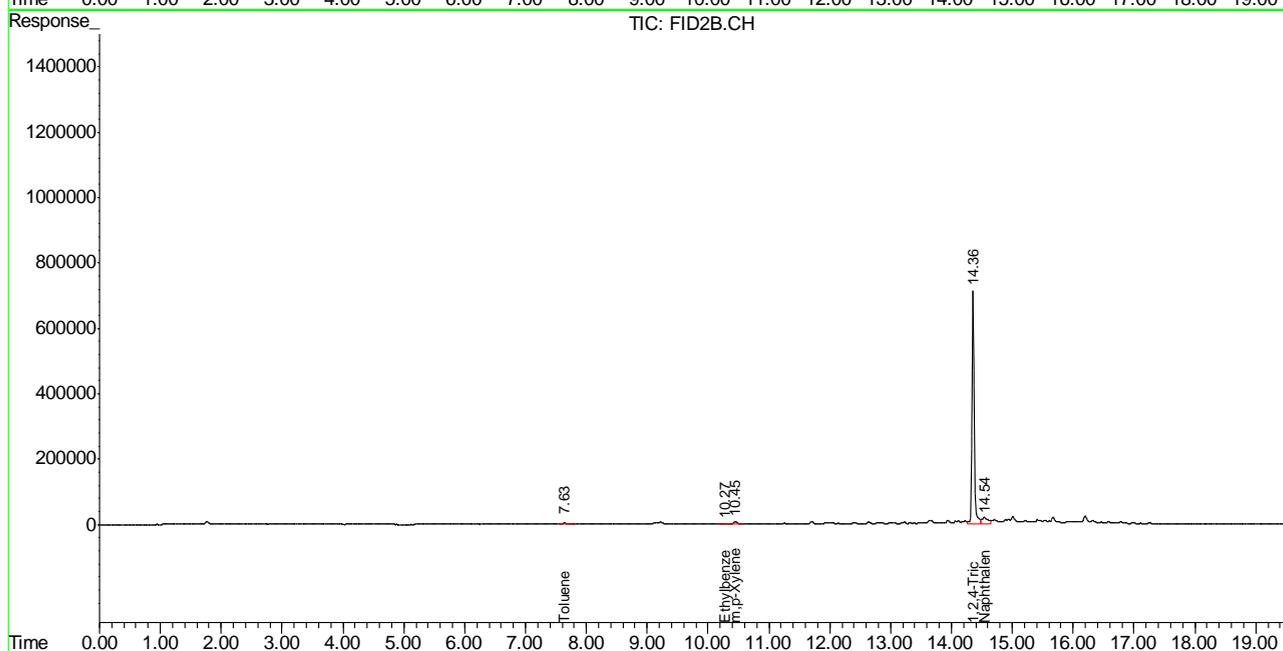
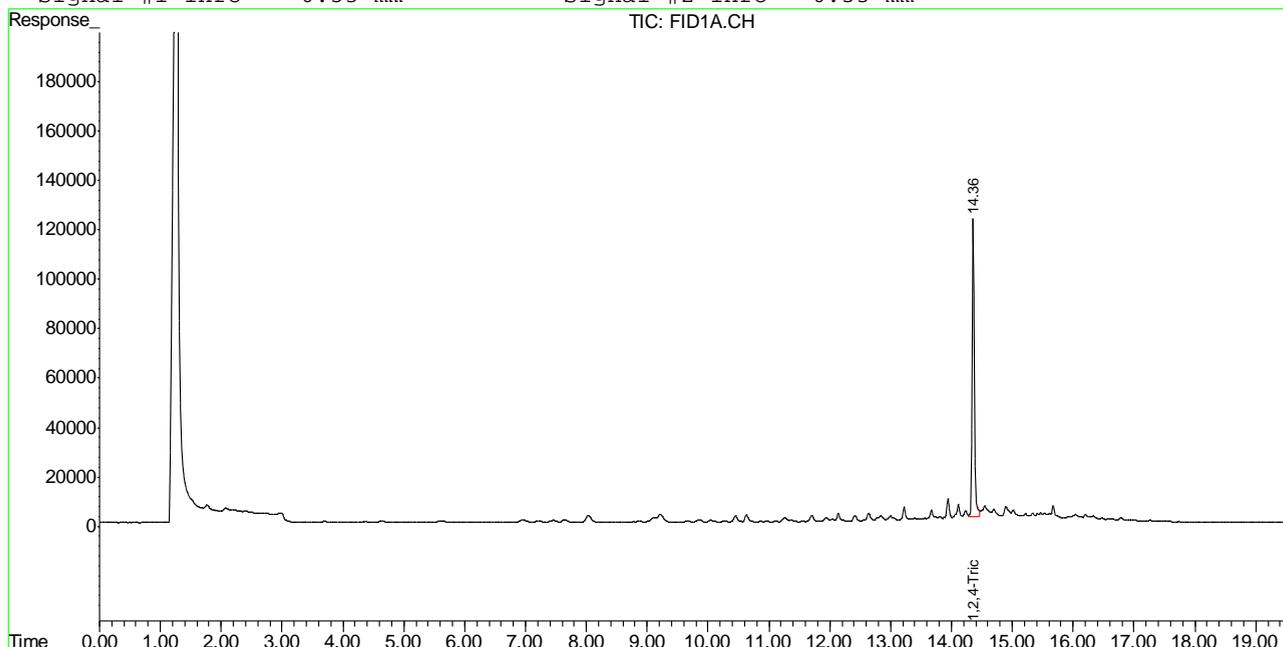
(f)=RT Delta > 1/2 Window (m)=manual int.  
 GB16420.D TB868GB868SOIL.M Fri Jun 22 08:12:58 2012 GC

Quantitation Report (QT Reviewed)

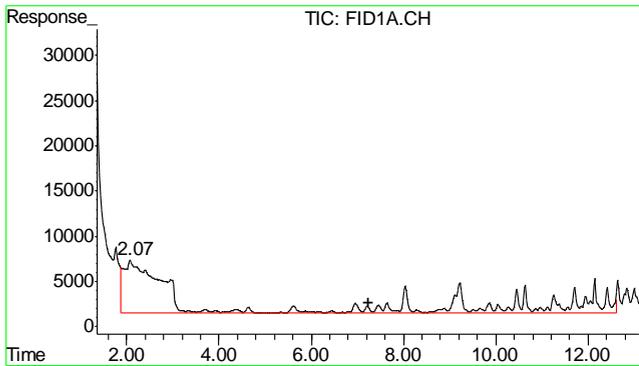
Signal #1 : Y:\1\DATA\062112\GB16420.D\FID1A.CH Vial: 9  
 Signal #2 : Y:\1\DATA\062112\GB16420.D\FID2B.CH  
 Acq On : 21 Jun 2012 7:05 pm Operator: StephK  
 Sample : D35710-1, 50X Inst : GC/MS Ins  
 Misc : GC2928,GGB910,5.044,,100,5,1 Multiplr: 1.00  
 IntFile Signal #1: TVH1.E IntFile Signal #2: FB2.E  
 Quant Time: Jun 22 7:13 2012 Quant Results File: TB868GB868SOIL.RES

Quant Method : C:\MSDCHEM\1...\TB868GB868SOIL.M (Chemstation Integrator)  
 Title : 8015B/8021B TVH/BTEX  
 Last Update : Thu Jun 21 15:17:12 2012  
 Response via : Multiple Level Calibration  
 DataAcq Meth : TVB4.M

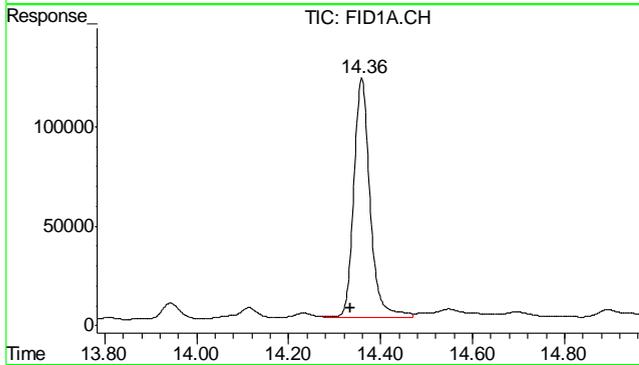
Volume Inj. :  
 Signal #1 Phase : DB-624 Signal #2 Phase: DB-624  
 Signal #1 Info : 0.53 mm Signal #2 Info : 0.53 mm



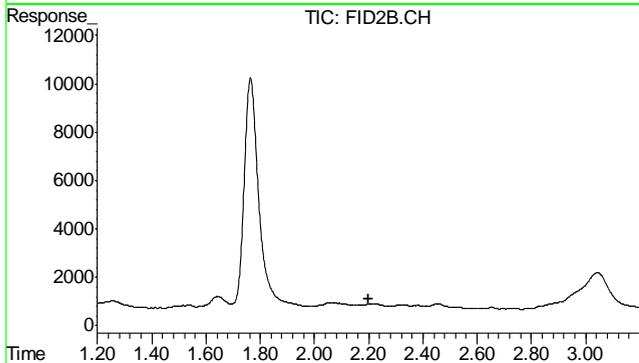
10.1.1  
10



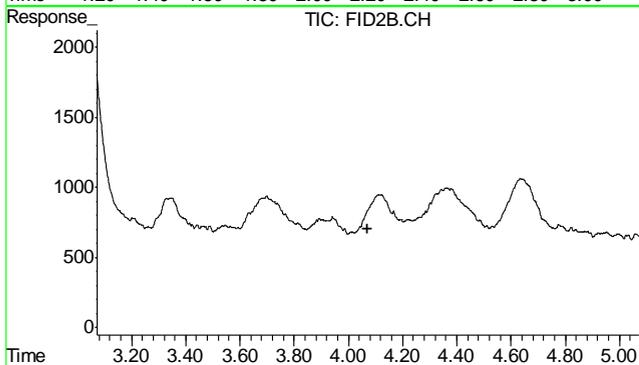
#1 TVH-Gasoline  
 R.T.: 7.230 min  
 Delta R.T.: 0.000 min  
 Response: 5980887  
 Conc: N.D.



#2 1,2,4-Trichlorobenzene  
 R.T.: 14.358 min  
 Delta R.T.: 0.024 min  
 Response: 2982648  
 Conc: 95.19 % m

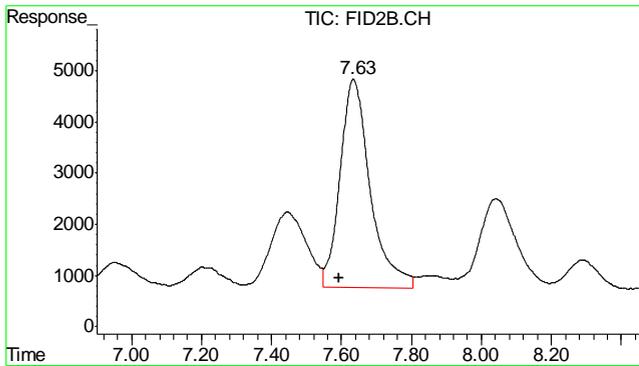


#4 Methyl-t-butyl-ether  
 R.T.: 0.000 min  
 Exp R.T.: 2.198 min  
 Response: 0  
 Conc: N.D.

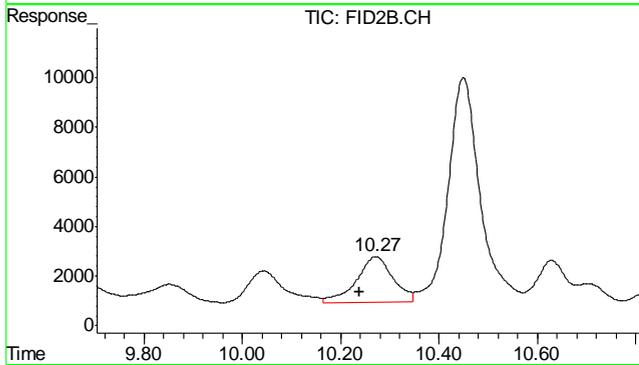


#5 Benzene  
 R.T.: 0.000 min  
 Exp R.T.: 4.072 min  
 Response: 0  
 Conc: N.D.

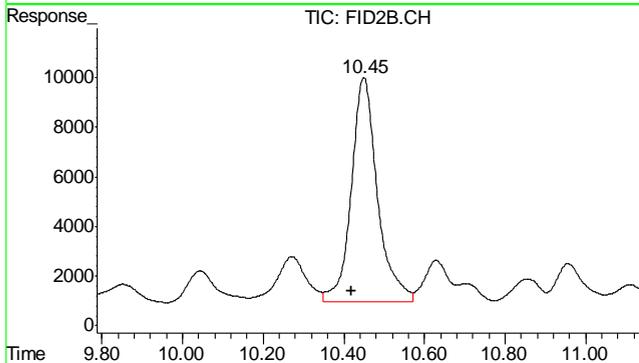
10.1.1  
 10



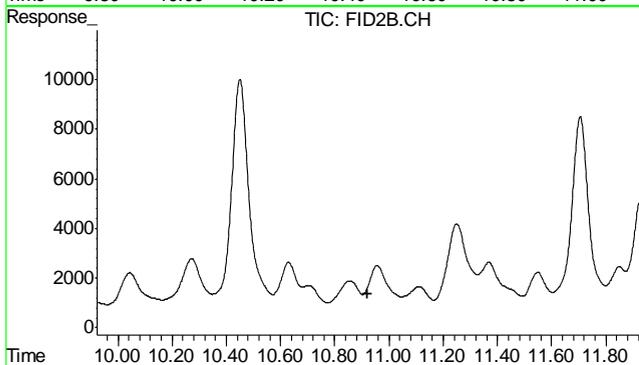
#6 Toluene  
 R.T.: 7.634 min  
 Delta R.T.: 0.040 min  
 Response: 244379  
 Conc: 0.62 ug/L



#7 Ethylbenzene  
 R.T.: 10.272 min  
 Delta R.T.: 0.034 min  
 Response: 95381  
 Conc: 0.28 ug/L

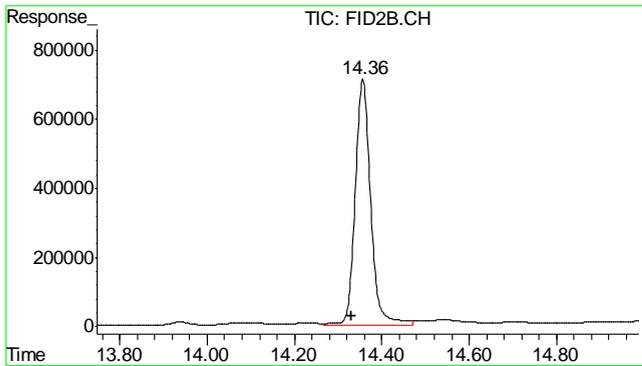


#8 m,p-Xylene  
 R.T.: 10.450 min  
 Delta R.T.: 0.029 min  
 Response: 407218  
 Conc: 0.74 ug/L



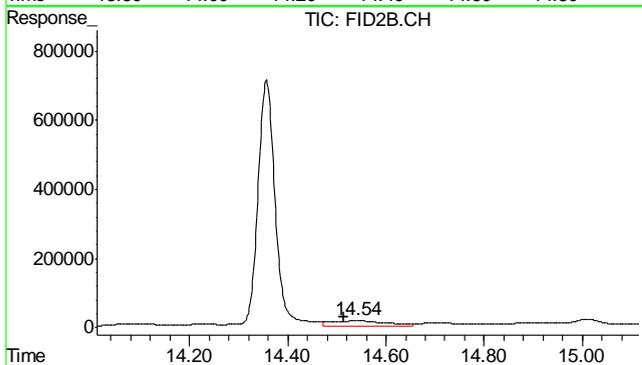
#9 o-Xylene  
 R.T.: 0.000 min  
 Exp R.T.: 10.922 min  
 Response: 0  
 Conc: N.D.

10.1.1  
**10**



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

R.T.: 14.357 min  
 Delta R.T.: 0.025 min  
 Response: 17634173  
 Conc: 108.50 %



#11 Naphthalene

R.T.: 14.541 min  
 Delta R.T.: 0.027 min  
 Response: 1354212  
 Conc: 6.86 ug/L

10.1.1  
 10

Quantitation Report (QT Reviewed)

Signal #1 : Y:\1\DATA\062112\GB16414.D\FID1A.CH Vial: 3  
 Signal #2 : Y:\1\DATA\062112\GB16414.D\FID2B.CH  
 Acq On : 21 Jun 2012 3:32 pm Operator: StephK  
 Sample : MB Inst : GC/MS Ins  
 Misc : GC2928,GGB910,5.000,,100,5,1 Multiplr: 1.00  
 IntFile Signal #1: TVH1.E IntFile Signal #2: FB2.E  
 Quant Time: Jun 21 15:56:52 2012 Quant Results File: TB868GB868SOIL.RES

Quant Method : C:\MSDCHEM\1...\TB868GB868SOIL.M (Chemstation Integrator)  
 Title : 8015B/8021B TVH/BTEX  
 Last Update : Thu Jun 21 15:17:12 2012  
 Response via : Initial Calibration  
 DataAcq Meth : TVB4.M

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

Compound	R.T.	Response	Conc	Units
-----				
System Monitoring Compounds				
2) S 1,2,4-Trichlorobenzene	14.35	3123776	99.693	%
10) S 1,2,4-Trichlorobenzene (P)	14.35	17660046	108.659	%
Target Compounds				
1) H TVH-Gasoline	7.23	4200171	<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	7.62	142182	0.359	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	14.52	245477	1.244	ug/L

(f)=RT Delta > 1/2 Window (m)=manual int.  
 GB16414.D TB868GB868SOIL.M Fri Jun 22 08:12:40 2012 GC

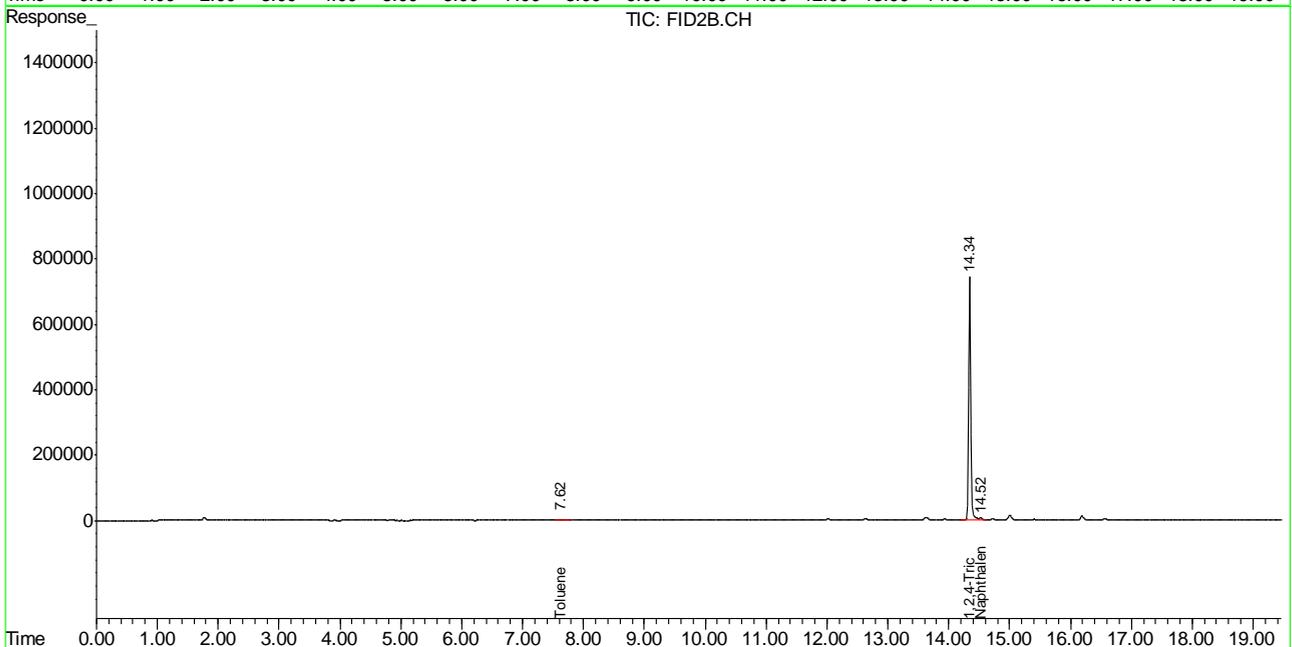
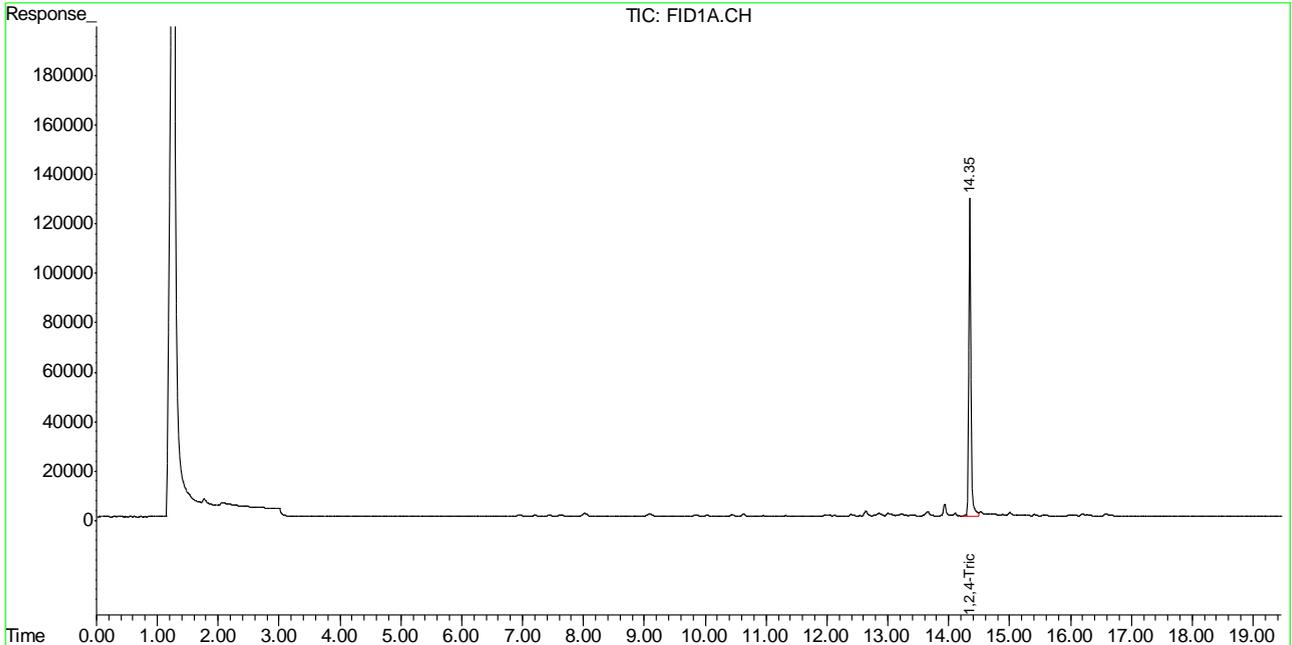
10.2.1  
 10

Quantitation Report (QT Reviewed)

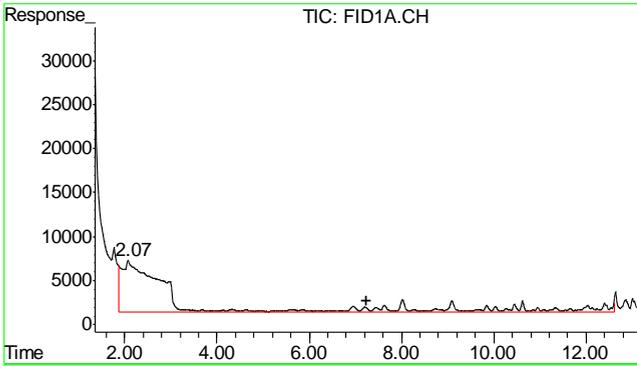
Signal #1 : Y:\1\DATA\062112\GB16414.D\FID1A.CH Vial: 3  
 Signal #2 : Y:\1\DATA\062112\GB16414.D\FID2B.CH  
 Acq On : 21 Jun 2012 3:32 pm Operator: StephK  
 Sample : MB Inst : GC/MS Ins  
 Misc : GC2928,GGB910,5.000,,100,5,1 Multiplr: 1.00  
 IntFile Signal #1: TVH1.E IntFile Signal #2: FB2.E  
 Quant Time: Jun 21 14:59 2012 Quant Results File: TB868GB868SOIL.RES

Quant Method : C:\MSDCHEM\1...\TB868GB868SOIL.M (Chemstation Integrator)  
 Title : 8015B/8021B TVH/BTEX  
 Last Update : Thu Jun 21 15:17:12 2012  
 Response via : Multiple Level Calibration  
 DataAcq Meth : TVB4.M

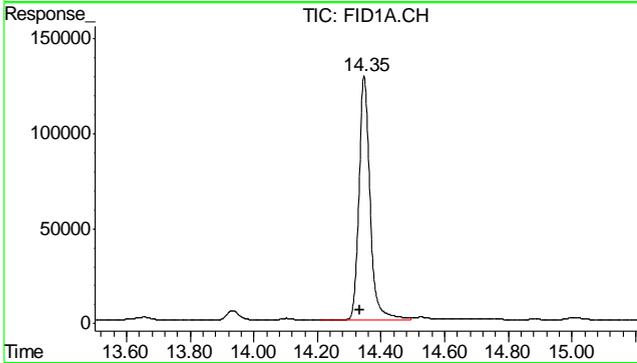
Volume Inj. :  
 Signal #1 Phase : DB-624 Signal #2 Phase: DB-624  
 Signal #1 Info : 0.53 mm Signal #2 Info : 0.53 mm



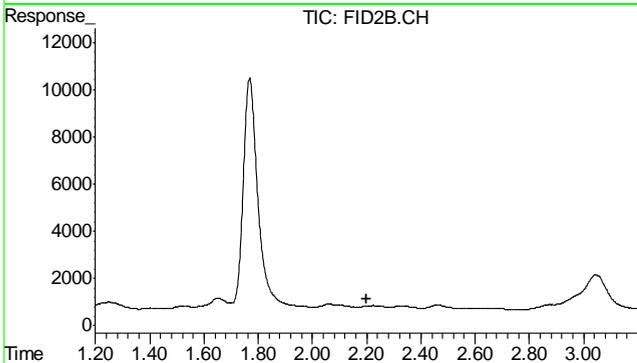
10.2.1  
10



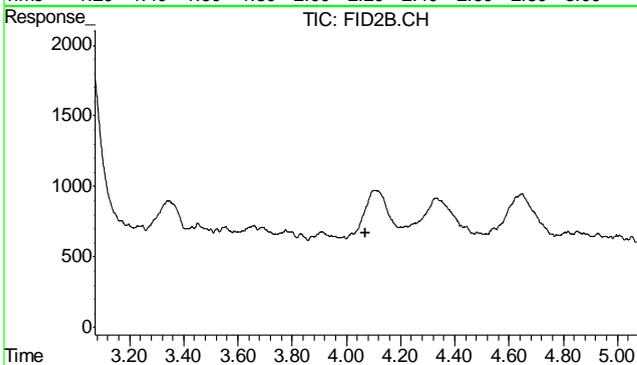
#1 TVH-Gasoline  
 R.T.: 7.230 min  
 Delta R.T.: 0.000 min  
 Response: 4200171  
 Conc: N.D.



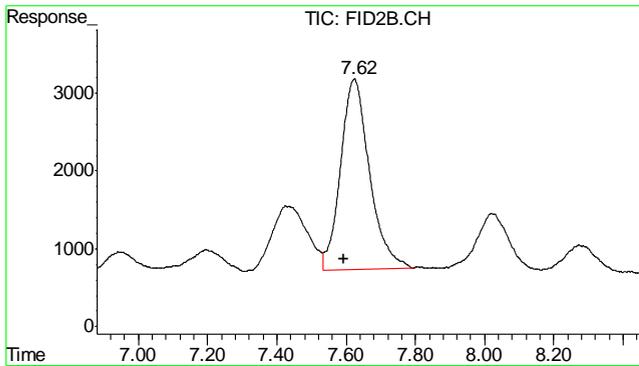
#2 1,2,4-Trichlorobenzene  
 R.T.: 14.347 min  
 Delta R.T.: 0.013 min  
 Response: 3123776  
 Conc: 99.69 %



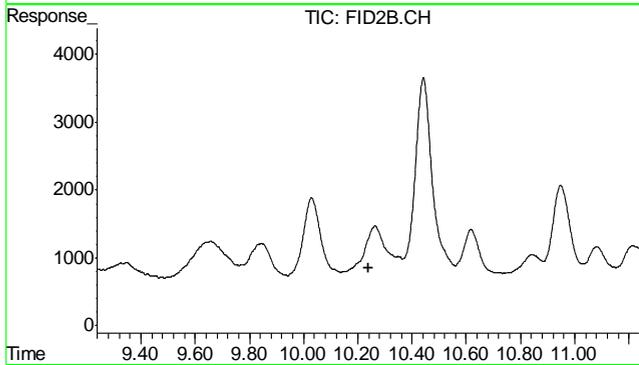
#4 Methyl-t-butyl-ether  
 R.T.: 0.000 min  
 Exp R.T.: 2.198 min  
 Response: 0  
 Conc: N.D.



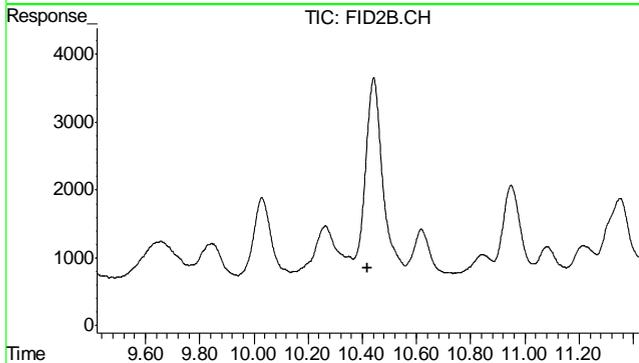
#5 Benzene  
 R.T.: 0.000 min  
 Exp R.T.: 4.072 min  
 Response: 0  
 Conc: N.D.



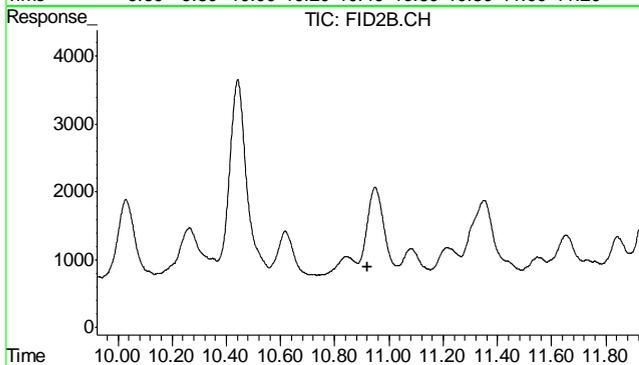
#6 Toluene  
 R.T.: 7.624 min  
 Delta R.T.: 0.030 min  
 Response: 142182  
 Conc: 0.36 ug/L



#7 Ethylbenzene  
 R.T.: 0.000 min  
 Exp R.T. : 10.238 min  
 Response: 0  
 Conc: N.D.

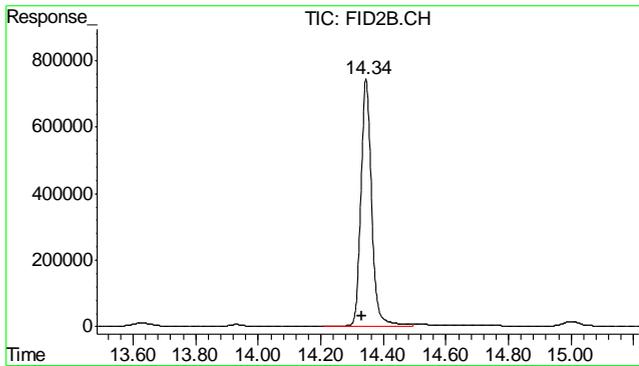


#8 m,p-Xylene  
 R.T.: 0.000 min  
 Exp R.T. : 10.421 min  
 Response: 0  
 Conc: N.D.



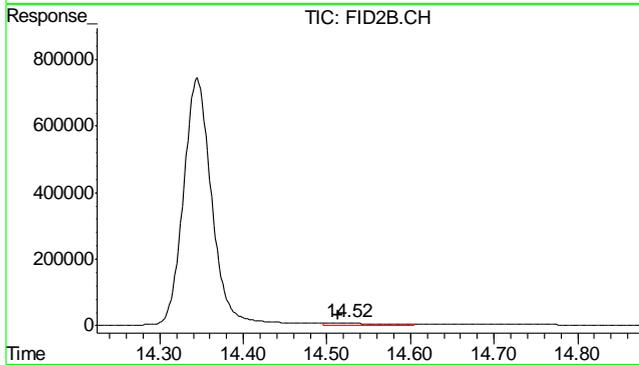
#9 o-Xylene  
 R.T.: 0.000 min  
 Exp R.T. : 10.922 min  
 Response: 0  
 Conc: N.D.

10.2.1  
 10



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

R.T.: 14.345 min  
Delta R.T.: 0.013 min  
Response: 17660046  
Conc: 108.66 %



#11 Naphthalene

R.T.: 14.524 min  
Delta R.T.: 0.011 min  
Response: 245477  
Conc: 1.24 ug/L

10.2.1  
10

## GC Semi-volatiles

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### QC Data Summaries

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Includes the following where applicable:

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



**Method Blank Summary**

**Job Number:** D35710  
**Account:** XTOKRWR XTO Energy  
**Project:** FRU 297-8B

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
OP6112-MB	FD14698.D	1	06/23/12	AW	06/22/12	OP6112	GFD764

The QC reported here applies to the following samples:

Method: SW846-8015B

D35710-1

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

CAS No.	Surrogate Recoveries	Limits
84-15-1	o-Terphenyl	96% 43-136%

11.1.1  
11

# Blank Spike Summary

**Job Number:** D35710  
**Account:** XTOKRWR XTO Energy  
**Project:** FRU 297-8B

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
OP6112-BS	FD14700.D	1	06/23/12	AW	06/22/12	OP6112	GFD764

The QC reported here applies to the following samples:

Method: SW846-8015B

D35710-1

CAS No.	Compound	Spike mg/kg	BSP mg/kg	BSP %	Limits
	TPH-DRO (C10-C28)	667	500	75	58-130

CAS No.	Surrogate Recoveries	BSP	Limits
84-15-1	o-Terphenyl	89%	43-136%

\* = Outside of Control Limits.

# Matrix Spike/Matrix Spike Duplicate Summary

**Job Number:** D35710  
**Account:** XTOKRWR XTO Energy  
**Project:** FRU 297-8B

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
OP6112-MS	FD14702.D	1	06/23/12	AW	06/22/12	OP6112	GFD764
OP6112-MSD	FD14704.D	1	06/23/12	AW	06/22/12	OP6112	GFD764
D35708-1	FD14706.D	1	06/23/12	AW	06/22/12	OP6112	GFD764

The QC reported here applies to the following samples:

Method: SW846-8015B

D35710-1

CAS No.	Compound	D35708-1 mg/kg	Spike mg/kg	MS mg/kg	MS %	MSD mg/kg	MSD %	RPD	Limits Rec/RPD
	TPH-DRO (C10-C28)	107	702	513	58	496	56	3	20-183/43

CAS No.	Surrogate Recoveries	MS	MSD	D35708-1	Limits
84-15-1	o-Terphenyl	77%	90%	96%	43-136%

11.3.1  
11

\* = Outside of Control Limits.

GC Semi-volatiles

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Raw Data

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Quantitation Report (QT Reviewed)

Data File : C:\MSDCHEM\2\DATA\2012\JUNE\FD062312.SEC\FD14720.D Vial: 64  
Acq On : 6-23-2012 05:19:31 PM Operator: alexwl  
Sample : D35710-1 Inst : FID5  
Misc : OP6112,GFD764,30.04,,,2,1 Multiplr: 1.00  
IntFile : autoint1.e  
Quant Time: Jun 25 09:38:51 2012 Quant Results File: DRO-GFD743R.RES

Quant Method : C:\MSDCHEM\2...\DRO-GFD743R.M (Chemstation Integrator)  
Title : 8015B TEH  
Last Update : Tue Jun 12 11:16:41 2012  
Response via : Initial Calibration  
DataAcq Meth : DRODUAL.M

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

Compound	R.T.	Response	Conc Units
-----			
System Monitoring Compounds			
1) S O-Terphenyl	9.63	35891973	826.693 mg/L
Target Compounds			
2) H TPH-DRO (c10-c28)	7.40	48196600	1160.651 mg/L

12.1.1  
**12**

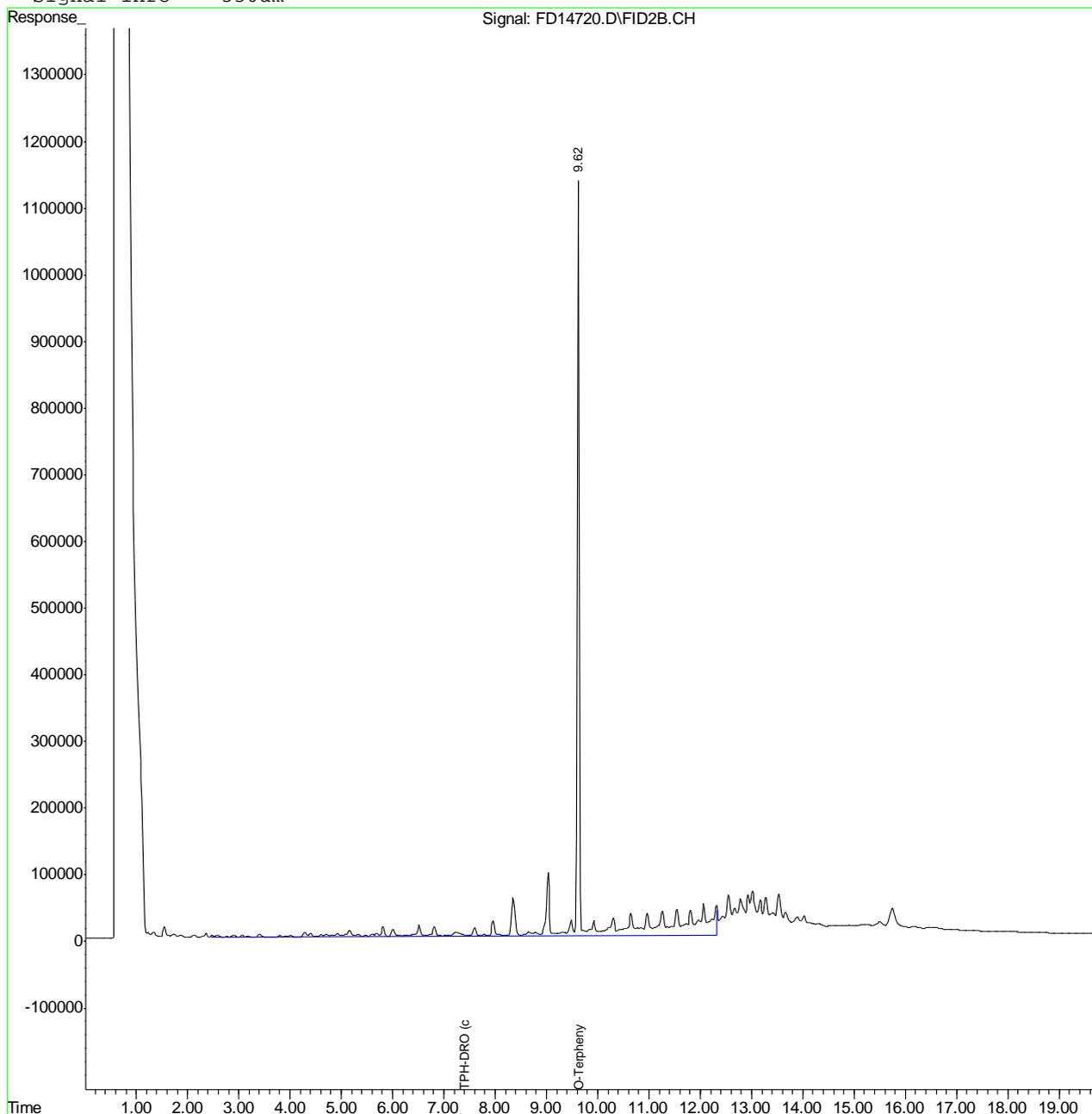
-----  
(f)=RT Delta > 1/2 Window (m)=manual int.  
FD14720.D DRO-GFD743R.M Mon Jun 25 10:37:27 2012 GC

Quantitation Report (QT Reviewed)

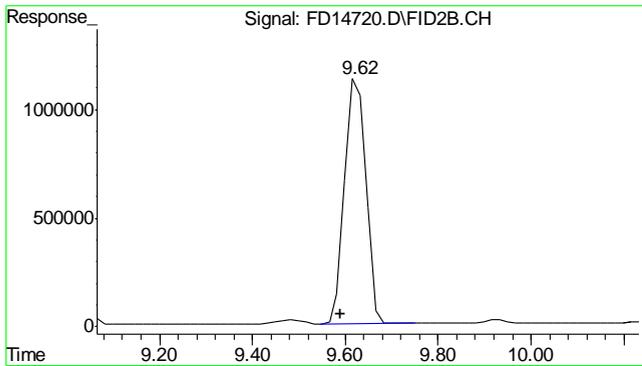
Data File : C:\MSDCHEM\2\DATA\2012\JUNE\FD062312.SEC\FD14720.D Vial: 64  
 Acq On : 6-23-2012 05:19:31 PM Operator: alexwl  
 Sample : D35710-1 Inst : FID5  
 Misc : OP6112,GFD764,30.04,,,2,1 Multiplr: 1.00  
 IntFile : autoint1.e  
 Quant Time: Jun 25 9:38 2012 Quant Results File: DRO-GFD743R.RES

Quant Method : C:\MSDCHEM\2...\DRO-GFD743R.M (Chemstation Integrator)  
 Title : 8015B TEH  
 Last Update : Tue Jun 12 11:16:41 2012  
 Response via : Multiple Level Calibration  
 DataAcq Meth : DRODUAL.M

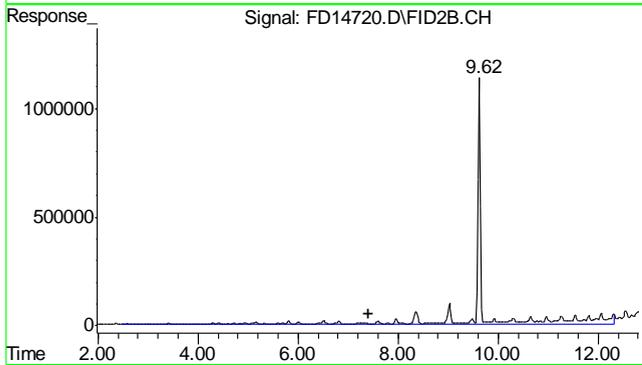
Volume Inj. : 1ul  
 Signal Phase : RTX-5  
 Signal Info : 530um



12.1.1  
**12**



#1 O-Terphenyl  
R.T.: 9.629 min  
Delta R.T.: 0.039 min  
Response: 35891973  
Conc: 826.69 mg/L



#2 TPH-DRO (c10-c28)  
R.T.: 7.400 min  
Delta R.T.: 0.000 min  
Response: 48196600  
Conc: 1160.65 mg/L m

12.1.1  
12

Quantitation Report (QT Reviewed)

Data File : C:\MSDCHEM\2\DATA\2012\JUNE\FD062312.SEC\FD14698.D Vial: 53  
 Acq On : 23 Jun 2012 12:31 pm Operator: alexwl  
 Sample : OP6112-MB Inst : FID5  
 Misc : OP6112,GFD764,30.00,,,2,1 Multiplr: 1.00  
 IntFile : autoint1.e  
 Quant Time: Jun 25 09:33:29 2012 Quant Results File: DRO-GFD743R.RES

Quant Method : C:\MSDCHEM\2...\DRO-GFD743R.M (Chemstation Integrator)  
 Title : 8015B TEH  
 Last Update : Tue Jun 12 11:16:41 2012  
 Response via : Initial Calibration  
 DataAcq Meth : DRODUAL.M

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

Compound	R.T.	Response	Conc Units
-----			
System Monitoring Compounds			
1) S O-Terphenyl	9.62	41812248	963.054 mg/L
Target Compounds			
2) H TPH-DRO (c10-c28)	7.40	2168150	52.213 mg/L

12.2.1  
**12**

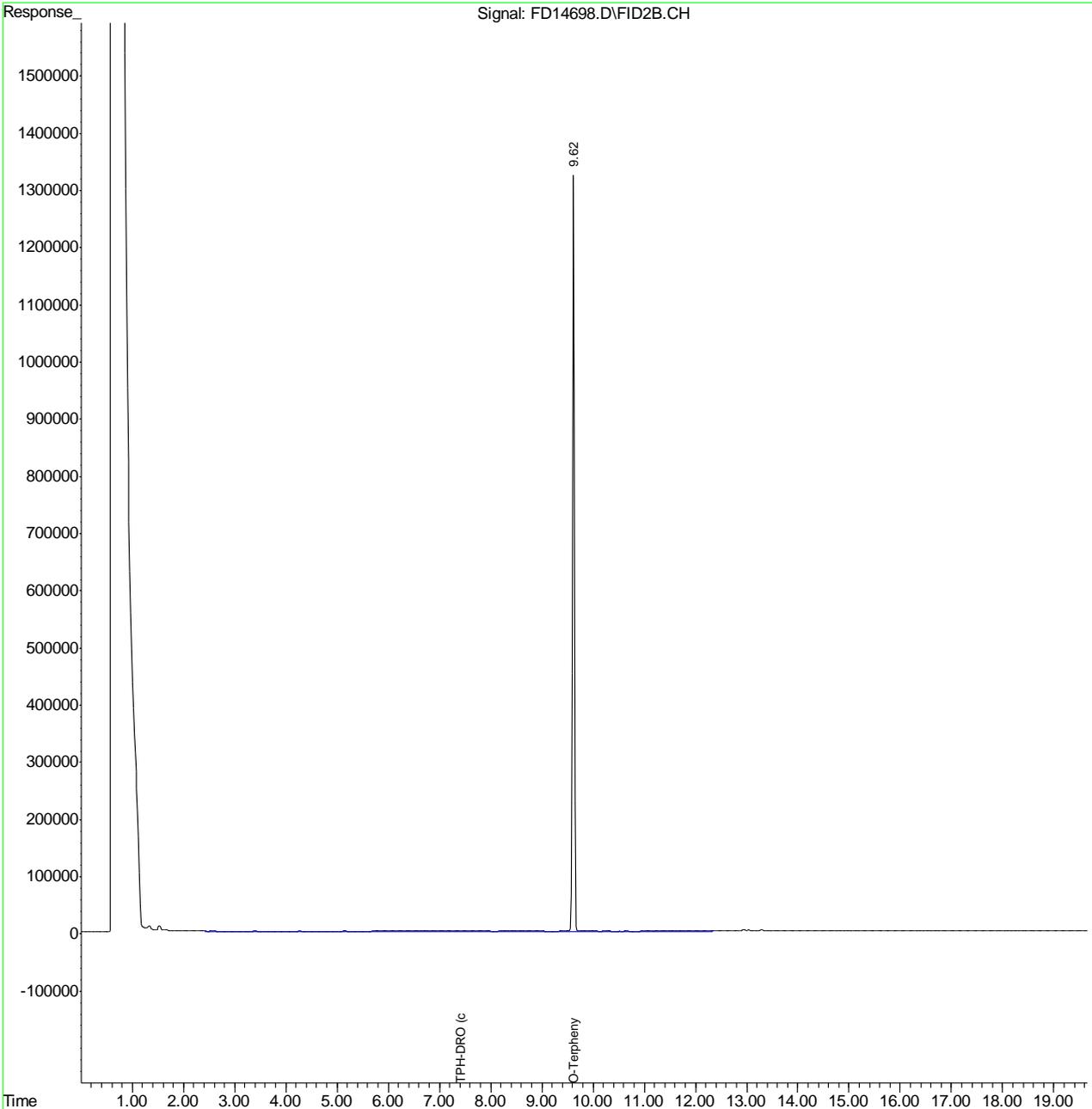
-----  
 (f)=RT Delta > 1/2 Window (m)=manual int.  
 FD14698.D DRO-GFD743R.M Mon Jun 25 10:37:19 2012 GC

Quantitation Report (QT Reviewed)

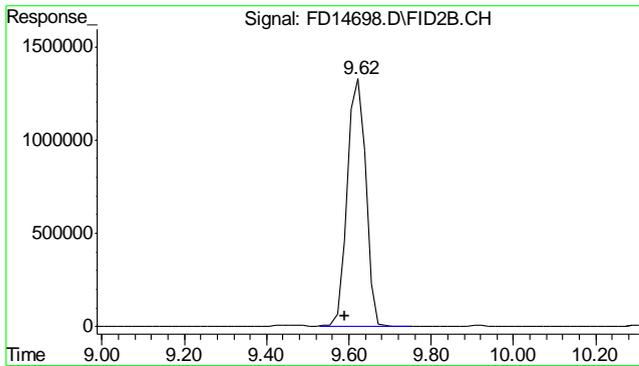
Data File : C:\MSDCHEM\2\DATA\2012\JUNE\FD062312.SEC\FD14698.D Vial: 53  
Acq On : 23 Jun 2012 12:31 pm Operator: alexwl  
Sample : OP6112-MB Inst : FID5  
Misc : OP6112,GFD764,30.00,,,2,1 Multiplr: 1.00  
IntFile : autoint1.e  
Quant Time: Jun 25 9:34 2012 Quant Results File: DRO-GFD743R.RES

Quant Method : C:\MSDCHEM\2...\DRO-GFD743R.M (Chemstation Integrator)  
Title : 8015B TEH  
Last Update : Tue Jun 12 11:16:41 2012  
Response via : Multiple Level Calibration  
DataAcq Meth : DRODUAL.M

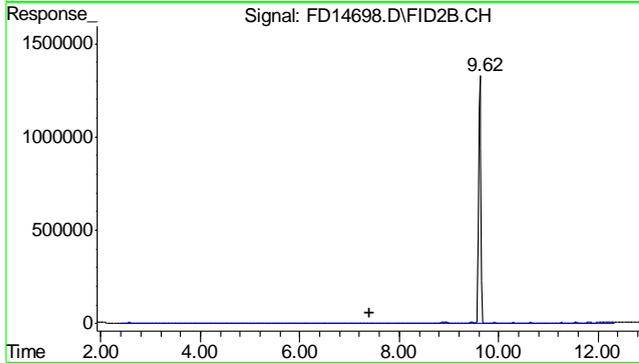
Volume Inj. : 1ul  
Signal Phase : RTX-5  
Signal Info : 530um



12.2.1  
**12**



#1 O-Terphenyl  
R.T.: 9.625 min  
Delta R.T.: 0.035 min  
Response: 41812248  
Conc: 963.05 mg/L



#2 TPH-DRO (c10-c28)  
R.T.: 7.400 min  
Delta R.T.: 0.000 min  
Response: 2168150  
Conc: 52.21 mg/L m

12.2.1  
**12**

## Metals Analysis

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### QC Data Summaries

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Includes the following where applicable:

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

BLANK RESULTS SUMMARY  
Part 2 - Method Blanks

Login Number: D35710  
Account: XTOKRWR - XTO Energy  
Project: FRU 297-8B

QC Batch ID: MP7737  
Matrix Type: SOLID

Methods: SW846 6010C  
Units: mg/kg

Prep Date: 06/22/12

Metal	RL	IDL	MDL	MB raw	final
Aluminum	10	.96	.57		
Antimony	3.0	.17	.12		
Arsenic	2.5	.44	.56		
Barium	1.0	.01	.11	0.060	<1.0
Beryllium	1.0	.13	.15		
Boron	5.0	.1	.06		
Cadmium	1.0	.06	.036	0.040	<1.0
Calcium	40	.54	9		
Chromium	1.0	.03	.03	0.070	<1.0
Cobalt	0.50	.04	.07		
Copper	1.0	.12	.15	0.15	<1.0
Iron	7.0	.12	.87		
Lead	5.0	.19	.24	-0.070	<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	-0.060	<3.0
Phosphorus	10	1.4	1.9		
Potassium	200	6.1	7		
Selenium	5.0	.48	.36	0.0	<5.0
Silicon	5.0	.29	.37		
Silver	3.0	.04	.06	-0.060	<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	-0.31	<3.0

Associated samples MP7737: D35710-1

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

13.11  
13

BLANK RESULTS SUMMARY  
Part 2 - Method Blanks

Login Number: D35710  
Account: XTOKRWR - XTO Energy  
Project: FRU 297-8B

QC Batch ID: MP7737  
Matrix Type: SOLID

Methods: SW846 6010C  
Units: mg/kg

Prep Date:

Metal

(anr) Analyte not requested

13.1.1  
13

MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D35710  
 Account: XTOKRWR - XTO Energy  
 Project: FRU 297-8B

QC Batch ID: MP7737  
 Matrix Type: SOLID

Methods: SW846 6010C  
 Units: mg/kg

Prep Date: 06/22/12

Metal	D35708-1 Original MS		SpikeLot ICPALL2	% Rec	QC Limits
Aluminum					
Antimony					
Arsenic					
Barium	807	954	213	69.1N(a)	75-125
Beryllium					
Boron					
Cadmium	0.40	47.6	53.2	88.8	75-125
Calcium					
Chromium	36.5	83.6	53.2	88.6	75-125
Cobalt					
Copper	11.3	61.2	53.2	93.9	75-125
Iron					
Lead	11.5	103	106	86.1	75-125
Lithium					
Magnesium					
Manganese					
Molybdenum					
Nickel	15.4	60.2	53.2	84.3	75-125
Phosphorus					
Potassium					
Selenium	1.4	95.4	106	88.4	75-125
Silicon					
Silver	0.0	17.6	21.3	82.8	75-125
Sodium					
Strontium					
Thallium					
Tin					
Titanium					
Uranium					
Vanadium					
Zinc	44.2	89.3	53.2	84.8	75-125

Associated samples MP7737: D35710-1

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

13.12  
13

MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D35710  
Account: XTOKRWR - XTO Energy  
Project: FRU 297-8B

QC Batch ID: MP7737  
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 recovery indicates possible matrix interference.

MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D35710  
 Account: XTOKRWR - XTO Energy  
 Project: FRU 297-8B

QC Batch ID: MP7737  
 Matrix Type: SOLID

Methods: SW846 6010C  
 Units: mg/kg

Prep Date: 06/22/12

Metal	D35708-1 Original	MSD	SpikeLot ICPAL2	% Rec	MSD RPD	QC Limit
Aluminum						
Antimony						
Arsenic						
Barium	807	996	215	88.0	4.3	20
Beryllium						
Boron						
Cadmium	0.40	47.7	53.7	88.1	0.2	20
Calcium						
Chromium	36.5	83.6	53.7	87.7	0.0	20
Cobalt						
Copper	11.3	61.5	53.7	93.5	0.5	20
Iron						
Lead	11.5	102	107	84.3	1.0	20
Lithium						
Magnesium						
Manganese						
Molybdenum						
Nickel	15.4	60.2	53.7	83.4	0.0	20
Phosphorus						
Potassium						
Selenium	1.4	95.8	107	87.9	0.4	20
Silicon						
Silver	0.0	17.8	21.5	82.9	1.1	20
Sodium						
Strontium						
Thallium						
Tin						
Titanium						
Uranium						
Vanadium						
Zinc	44.2	88.4	53.7	82.3	1.0	20

Associated samples MP7737: D35710-1

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

13.12  
13

MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D35710  
Account: XTOKRWR - XTO Energy  
Project: FRU 297-8B

QC Batch ID: MP7737  
Matrix Type: SOLID

Methods: SW846 6010C  
Units: mg/kg

Prep Date:

Metal

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

SPIKE BLANK AND LAB CONTROL SAMPLE SUMMARY

Login Number: D35710  
 Account: XTOKRWR - XTO Energy  
 Project: FRU 297-8B

QC Batch ID: MP7737  
 Matrix Type: SOLID

Methods: SW846 6010C  
 Units: mg/kg

Prep Date: 06/22/12

Metal	BSP Result	Spikelot ICPALL2	% Rec	QC Limits
Aluminum				
Antimony				
Arsenic				
Barium	207	200	103.5	80-120
Beryllium				
Boron				
Cadmium	49.5	50	99.0	80-120
Calcium				
Chromium	52.7	50	105.4	80-120
Cobalt				
Copper	49.1	50	98.2	80-120
Iron				
Lead	99.3	100	99.3	80-120
Lithium				
Magnesium				
Manganese				
Molybdenum				
Nickel	50.2	50	100.4	80-120
Phosphorus				
Potassium				
Selenium	98.6	100	98.6	80-120
Silicon				
Silver	20.3	20	101.5	80-120
Sodium				
Strontium				
Thallium				
Tin				
Titanium				
Uranium				
Vanadium				
Zinc	50.6	50	101.2	80-120

Associated samples MP7737: D35710-1

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

13.1.3  
 13

SPIKE BLANK AND LAB CONTROL SAMPLE SUMMARY

Login Number: D35710  
Account: XTOKRWR - XTO Energy  
Project: FRU 297-8B

QC Batch ID: MP7737  
Matrix Type: SOLID

Methods: SW846 6010C  
Units: mg/kg

Prep Date:

Metal

(anr) Analyte not requested

SERIAL DILUTION RESULTS SUMMARY

Login Number: D35710  
 Account: XTOKRWR - XTO Energy  
 Project: FRU 297-8B

QC Batch ID: MP7737  
 Matrix Type: SOLID

Methods: SW846 6010C  
 Units: ug/l

Prep Date: 06/22/12

Metal	D35708-1 Original SDL 1:5		%DIF	QC Limits
Aluminum				
Antimony				
Arsenic				
Barium	7890	8870	12.4*(a)	0-10
Beryllium				
Boron				
Cadmium	3.90	0.00	100.0(b)	0-10
Calcium				
Chromium	358	407	13.8*(a)	0-10
Cobalt				
Copper	111	117	5.0	0-10
Iron				
Lead	113	114	0.4	0-10
Lithium				
Magnesium				
Manganese				
Molybdenum				
Nickel	150	172	14.4*(a)	0-10
Phosphorus				
Potassium				
Selenium	13.9	39.0	180.6(b)	0-10
Silicon				
Silver	0.00	0.00	NC	0-10
Sodium				
Strontium				
Thallium				
Tin				
Titanium				
Uranium				
Vanadium				
Zinc	432	499	15.5*(a)	0-10

Associated samples MP7737: D35710-1

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

13.14  
13

SERIAL DILUTION RESULTS SUMMARY

Login Number: D35710  
Account: XTOKRWR - XTO Energy  
Project: FRU 297-8B

QC Batch ID: MP7737  
Matrix Type: SOLID

Methods: SW846 6010C  
Units: ug/l

Prep Date:

Metal

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

BLANK RESULTS SUMMARY  
Part 2 - Method Blanks

Login Number: D35710  
Account: XTOKRWR - XTO Energy  
Project: FRU 297-8B

QC Batch ID: MP7738  
Matrix Type: SOLID

Methods: SW846 6020A  
Units: mg/kg

Prep Date: 06/22/12

Metal	RL	IDL	MDL	MB raw	final
Aluminum	25	.22	.31		
Antimony	0.20	.0018	.0075		
Arsenic	0.10	.042	.06	0.0034	<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 MP7738: D35710-1

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

13.21  
13

MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D35710  
 Account: XTOKRWR - XTO Energy  
 Project: FRU 297-8B

QC Batch ID: MP7738  
 Matrix Type: SOLID

Methods: SW846 6020A  
 Units: mg/kg

Prep Date: 06/22/12

Metal	D35708-1 Original MS		Spikelot ICPALL2	% Rec	QC Limits
Aluminum					
Antimony					
Arsenic	8.9	131	106	114.8	75-125
Barium					
Beryllium					
Boron					
Cadmium					
Calcium					
Chromium					
Cobalt					
Copper					
Iron					
Lead					
Magnesium					
Manganese					
Molybdenum					
Nickel					
Phosphorus					
Potassium					
Selenium					
Silver					
Sodium					
Strontium					
Thallium					
Tin					
Titanium					
Uranium					
Vanadium					
Zinc					

Associated samples MP7738: D35710-1

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

13.22  
13

MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D35710  
 Account: XTOKRWR - XTO Energy  
 Project: FRU 297-8B

QC Batch ID: MP7738  
 Matrix Type: SOLID

Methods: SW846 6020A  
 Units: mg/kg

Prep Date: 06/22/12

Metal	D35708-1 Original MSD		SpikeLot ICPALL2	% Rec	MSD RPD	QC Limit
Aluminum						
Antimony						
Arsenic	8.9	113	107	96.9	14.8	20
Barium						
Beryllium						
Boron						
Cadmium						
Calcium						
Chromium						
Cobalt						
Copper						
Iron						
Lead						
Magnesium						
Manganese						
Molybdenum						
Nickel						
Phosphorus						
Potassium						
Selenium						
Silver						
Sodium						
Strontium						
Thallium						
Tin						
Titanium						
Uranium						
Vanadium						
Zinc						

Associated samples MP7738: D35710-1

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

13.22  
13

SPIKE BLANK AND LAB CONTROL SAMPLE SUMMARY

Login Number: D35710  
 Account: XTOKRWR - XTO Energy  
 Project: FRU 297-8B

QC Batch ID: MP7738  
 Matrix Type: SOLID

Methods: SW846 6020A  
 Units: mg/kg

Prep Date: 06/22/12

Metal	BSP Result	Spikelot ICPALL2	% Rec	QC Limits
Aluminum				
Antimony				
Arsenic	106	100	106.0	80-120
Barium				
Beryllium				
Boron				
Cadmium				
Calcium				
Chromium				
Cobalt				
Copper				
Iron				
Lead				
Magnesium				
Manganese				
Molybdenum				
Nickel				
Phosphorus				
Potassium				
Selenium				
Silver				
Sodium				
Strontium				
Thallium				
Tin				
Titanium				
Uranium				
Vanadium				
Zinc				

Associated samples MP7738: D35710-1

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

13.23  
13

SERIAL DILUTION RESULTS SUMMARY

Login Number: D35710  
 Account: XTOKRWR - XTO Energy  
 Project: FRU 297-8B

QC Batch ID: MP7738  
 Matrix Type: SOLID

Methods: SW846 6020A  
 Units: ug/l

Prep Date: 06/22/12

Metal	D35708-1			QC
	Original	SDL 5:25	%DIF	Limits
Aluminum				
Antimony				
Arsenic	86.6	87.1	0.5	0-10
Barium				
Beryllium				
Boron				
Cadmium				
Calcium				
Chromium				
Cobalt				
Copper				
Iron				
Lead				
Magnesium				
Manganese				
Molybdenum				
Nickel				
Phosphorus				
Potassium				
Selenium				
Silver				
Sodium				
Strontium				
Thallium				
Tin				
Titanium				
Uranium				
Vanadium				
Zinc				

Associated samples MP7738: D35710-1

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

13.24  
 13

BLANK RESULTS SUMMARY  
Part 2 - Method Blanks

Login Number: D35710  
Account: XTOKRWR - XTO Energy  
Project: FRU 297-8B

QC Batch ID: MP7743  
Matrix Type: AQUEOUS

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

Prep Date: 06/22/12

Metal	RL	IDL	MDL	MB raw	final
Aluminum	500	48	110		
Antimony	150	8.5	16		
Arsenic	130	22	38		
Barium	50	.5	2.5		
Beryllium	50	6.5	16		
Boron	250	5	13		
Cadmium	50	3	3		
Calcium	2000	27	37	414	<2000
Chromium	50	1.5	2		
Cobalt	25	2	2		
Copper	50	6	15		
Iron	350	6	95		
Lead	250	9.5	15		
Lithium	10	2.5	3.3		
Magnesium	1000	33	55	31.5	<1000
Manganese	25	6	9		
Molybdenum	50	11	11		
Nickel	150	2.5	2.7		
Phosphorus	500	70	300		
Potassium	5000	310	310		
Selenium	250	24	29		
Silicon	250	15	11		
Silver	150	2	3.3		
Sodium	2000	30	490	-270	<2000
Strontium	25	.2	7.5		
Thallium	50	15	15		
Tin	250	60	120		
Titanium	50	.5	6		
Uranium	250	11	11		
Vanadium	50	1	2		
Zinc	150	2.5	7.5		

Associated samples MP7743: D35710-1A

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

13.3.1  
13

BLANK RESULTS SUMMARY  
Part 2 - Method Blanks

Login Number: D35710  
Account: XTOKRWR - XTO Energy  
Project: FRU 297-8B

QC Batch ID: MP7743  
Matrix Type: AQUEOUS

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

Prep Date:

Metal

(anr) Analyte not requested

13.3.1  
13

MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D35710  
 Account: XTOKRWR - XTO Energy  
 Project: FRU 297-8B

QC Batch ID: MP7743  
 Matrix Type: AQUEOUS

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

Prep Date: 06/22/12

Metal	D35708-1A Original MS	Spikelot ICPALL2	% Rec	QC Limits	
Aluminum					
Antimony					
Arsenic					
Barium					
Beryllium					
Boron					
Cadmium					
Calcium	139000	281000	125000	113.6	75-125
Chromium					
Cobalt					
Copper					
Iron					
Lead					
Lithium					
Magnesium	16000	140000	125000	99.2	75-125
Manganese					
Molybdenum					
Nickel					
Phosphorus					
Potassium					
Selenium					
Silicon					
Silver					
Sodium	1620000	1780000	125000	128.0(a)	75-125
Strontium					
Thallium					
Tin					
Titanium					
Uranium					
Vanadium					
Zinc					

Associated samples MP7743: D35710-1A

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

13.32  
13

MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D35710  
Account: XTOKRWR - XTO Energy  
Project: FRU 297-8B

QC Batch ID: MP7743  
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
- (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: D35710  
 Account: XTOKRWR - XTO Energy  
 Project: FRU 297-8B

QC Batch ID: MP7743  
 Matrix Type: AQUEOUS

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

Prep Date: 06/22/12

Metal	D35708-1A Original MSD	Spikelot ICPALL2	% Rec	MSD RPD	QC Limit	
Aluminum						
Antimony						
Arsenic						
Barium						
Beryllium						
Boron						
Cadmium						
Calcium	139000	294000	125000	124.0	4.5	20
Chromium						
Cobalt						
Copper						
Iron						
Lead						
Lithium						
Magnesium	16000	142000	125000	100.8	1.4	20
Manganese						
Molybdenum						
Nickel						
Phosphorus						
Potassium						
Selenium						
Silicon						
Silver						
Sodium	1620000	1960000	125000	272.0(a)	9.6	20
Strontium						
Thallium						
Tin						
Titanium						
Uranium						
Vanadium						
Zinc						

Associated samples MP7743: D35710-1A

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

13.32  
13

MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D35710  
Account: XTOKRWR - XTO Energy  
Project: FRU 297-8B

QC Batch ID: MP7743  
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
- (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: D35710  
 Account: XTOKRWR - XTO Energy  
 Project: FRU 297-8B

QC Batch ID: MP7743  
 Matrix Type: AQUEOUS

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

Prep Date: 06/22/12

Metal	BSP Result	Spikelot ICPALL2	% Rec	QC Limits
Aluminum				
Antimony				
Arsenic				
Barium				
Beryllium				
Boron				
Cadmium				
Calcium	139000	125000	111.2	80-120
Chromium				
Cobalt				
Copper				
Iron				
Lead				
Lithium				
Magnesium	124000	125000	99.2	80-120
Manganese				
Molybdenum				
Nickel				
Phosphorus				
Potassium				
Selenium				
Silicon				
Silver				
Sodium	124000	125000	99.2	80-120
Strontium				
Thallium				
Tin				
Titanium				
Uranium				
Vanadium				
Zinc				

Associated samples MP7743: D35710-1A

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

13.3.3  
 13

SPIKE BLANK AND LAB CONTROL SAMPLE SUMMARY

Login Number: D35710  
Account: XTOKRWR - XTO Energy  
Project: FRU 297-8B

QC Batch ID: MP7743  
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: D35710  
 Account: XTOKRWR - XTO Energy  
 Project: FRU 297-8B

QC Batch ID: MP7743  
 Matrix Type: AQUEOUS

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

Prep Date: 06/22/12

Metal	D35708-1A Original SDL 1:5		%DIF	QC Limits
Aluminum				
Antimony				
Arsenic				
Barium				
Beryllium				
Boron				
Cadmium				
Calcium	27800	27700	0.4	0-10
Chromium				
Cobalt				
Copper				
Iron				
Lead				
Lithium				
Magnesium	3200	3340	4.4	0-10
Manganese				
Molybdenum				
Nickel				
Phosphorus				
Potassium				
Selenium				
Silicon				
Silver				
Sodium	323000	337000	4.3	0-10
Strontium				
Thallium				
Tin				
Titanium				
Uranium				
Vanadium				
Zinc				

Associated samples MP7743: D35710-1A

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

13.3.4  
 13

SERIAL DILUTION RESULTS SUMMARY

Login Number: D35710  
Account: XTOKRWR - XTO Energy  
Project: FRU 297-8B

QC Batch ID: MP7743  
Matrix Type: AQUEOUS

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

Prep Date:

Metal

(anr) Analyte not requested

13.3.4

13

BLANK RESULTS SUMMARY  
Part 2 - Method Blanks

Login Number: D35710  
Account: XTOKRWR - XTO Energy  
Project: FRU 297-8B

QC Batch ID: MP7755  
Matrix Type: SOLID

Methods: SW846 7471B  
Units: mg/kg

Prep Date: 06/26/12

Metal	RL	IDL	MDL	MB	
				raw	final
Mercury	0.10	.0011	.0009	0.00037	<0.10

Associated samples MP7755: D35710-1

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

MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D35710  
 Account: XTOKRWR - XTO Energy  
 Project: FRU 297-8B

QC Batch ID: MP7755  
 Matrix Type: SOLID

Methods: SW846 7471B  
 Units: mg/kg

Prep Date: 06/26/12

Metal	D35710-1 Original MS	Spike HGWSR1	Spikelot % Rec	QC Limits
Mercury	0.016	0.43	0.434 95.3	75-125

Associated samples MP7755: D35710-1

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

MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D35710  
 Account: XTOKRWR - XTO Energy  
 Project: FRU 297-8B

QC Batch ID: MP7755  
 Matrix Type: SOLID

Methods: SW846 7471B  
 Units: mg/kg

Prep Date: 06/26/12

Metal	D35710-1 Original MSD	Spike HGWSR1	lot % Rec	MSD RPD	QC Limit
Mercury	0.016	0.42	0.426	94.9	2.4

Associated samples MP7755: D35710-1

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

SPIKE BLANK AND LAB CONTROL SAMPLE SUMMARY

Login Number: D35710  
Account: XTOKRWR - XTO Energy  
Project: FRU 297-8B

QC Batch ID: MP7755  
Matrix Type: SOLID

Methods: SW846 7471B  
Units: mg/kg

Prep Date: 06/26/12

Metal	BSP Result	Spikelot HGWSR1	% Rec	QC Limits
Mercury	0.34	0.4	85.0	80-120

Associated samples MP7755: D35710-1

Results < IDL are shown as zero for calculation purposes  
(\* ) Outside of QC limits  
(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: D35710  
Account: XTOKRWR - XTO Energy  
Project: FRU 297-8B

Analyte	Batch ID	RL	MB Result	Units	Spike Amount	BSP Result	BSP %Recov	QC Limits
Chromium, Hexavalent	GP7549/GN15579	1.0	0.0	mg/kg	261	252	96.6	80-120%
Specific Conductivity	GP7560/GN15563			umhos/cm	10009	9870	98.6	90-110%
pH	GN15564			su	8.00	8.00	100.0	99.3-100.7%

Associated Samples:  
Batch GN15564: D35710-1  
Batch GP7549: D35710-1  
Batch GP7560: D35710-1  
(\* ) Outside of QC limits

14.1  
14

DUPLICATE RESULTS SUMMARY  
GENERAL CHEMISTRY

Login Number: D35710  
Account: XTOKRWR - XTO Energy  
Project: FRU 297-8B

Analyte	Batch ID	QC Sample	Units	Original Result	DUP Result	RPD	QC Limits
Chromium, Hexavalent	GP7549/GN15579	D35573-1	mg/kg	0.0	0.0	0.0	0-20%
Redox Potential Vs H2	GN15599	D35794-1	mv	273	266	2.6	0-20%

Associated Samples:  
Batch GN15599: D35710-1  
Batch GP7549: D35710-1  
(\* ) Outside of QC limits

14.2  
14

MATRIX SPIKE RESULTS SUMMARY  
GENERAL CHEMISTRY

Login Number: D35710  
Account: XTOKRWR - XTO Energy  
Project: FRU 297-8B

Analyte	Batch ID	QC Sample	Units	Original Result	Spike Amount	MS Result	%Rec	QC Limits
Chromium, Hexavalent	GP7549/GN15579	D35573-1	mg/kg	0.0	40	36.8	92.0	75-125%

Associated Samples:

Batch GP7549: D35710-1

(\*) Outside of QC limits

(N) Matrix Spike Rec. outside of QC limits

14.3

14

MATRIX SPIKE DUPLICATE RESULTS SUMMARY  
GENERAL CHEMISTRY

Login Number: D35710  
Account: XTOKRWR - XTO Energy  
Project: FRU 297-8B

Analyte	Batch ID	QC Sample	Units	Original Result	Spike Amount	MSD Result	RPD	QC Limit
Chromium, Hexavalent	GP7549/GN15579	D35573-1	mg/kg	0.0	40	37.3	1.4	

Associated Samples:

Batch GP7549: D35710-1

(\*) Outside of QC limits

(N) Matrix Spike Rec. outside of QC limits

14.4  
14