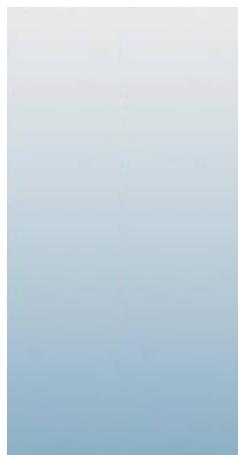




12/12/12



## Technical Report for

**XTO Energy**

**PCU 296-5A**

**1210-04**

**Accutest Job Number: D41506**

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

### Report to:

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ATTN: Dwayne Knudson**

**Total number of pages in report: 139**



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

A handwritten signature in black ink.

**Brad Madadian  
Laboratory Director**

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

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

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

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## Sample Summary

XTO Energy

**Job No:** D41506

PCU 296-5A

Project No: 1210-04

Sample Number	Collected Date	Time By	Matrix Received	Code Type	Client Sample ID
D41506-1	12/03/12	10:20 DS	12/05/12	SO Soil	RP SUBLINER(COMP)
D41506-1A	12/03/12	10:20 DS	12/05/12	SO Soil	RP SUBLINER(COMP)

---

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



## CASE NARRATIVE / CONFORMANCE SUMMARY

**Client:** XTO Energy

**Job No** D41506

**Site:** PCU 296-5A

**Report Date** 12/11/2012 2:52:51 PM

On 12/05/2012, 1 sample(s), 0 Trip Blank(s), and 0 Field Blank(s) were received at Accutest Mountain States (AMS) at a temperature of 3.2 °C. The samples were intact and properly preserved, unless noted below. An AMS Job Number of D41506 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> V3V1286
------------------	--------------------------

- All samples were analyzed within the recommended method holding time.
- Sample(s) D41506-1MS, D41506-1MSD were used as the QC samples indicated.
- All method blanks for this batch meet method specific criteria.

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

<b>Matrix</b> SO	<b>Batch ID:</b> OP7075
------------------	-------------------------

- All samples were extracted within the recommended method holding time.
- All samples were analyzed within the recommended method holding time.
- Sample(s) D41381-1MS, D41381-1MSD 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> GGB1023
------------------	--------------------------

- All samples were analyzed within the recommended method holding time.
- Sample(s) D41506-1MS, D41506-1MSD were used as the QC samples indicated.
- All method blanks for this batch meet method specific criteria.

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

<b>Matrix</b> SO	<b>Batch ID:</b> OP7061
------------------	-------------------------

- All samples were extracted within the recommended method holding time.
- All samples were analyzed within the recommended method holding time.
- Sample(s) D41506-1MS, D41506-1MSD 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:** MP9015

- All samples were digested within the recommended method holding time.
- All samples were analyzed within the recommended method holding time.
- All method blanks for this batch meet method specific criteria.
- Sample(s) D41381-1AMS, D41381-1AMSD, D41381-1ASDL were used as the QC samples for the metals analysis.

**Matrix** SO

**Batch ID:** MP9011

- All samples were digested within the recommended method holding time.
- All samples were analyzed within the recommended method holding time.
- All method blanks for this batch meet method specific criteria.
- Sample(s) D41506-1MS, D41506-1MSD, D41506-1SDL were used as the QC samples for the metals analysis.
- The matrix spike (MS) recovery(s) of Lead are outside control limits. Spike recovery indicates possible matrix interference.
- The matrix spike duplicate (MSD) recovery(s) of Lead, Nickel, Zinc are outside control limits. Probable cause due to matrix interference.
- The matrix spike (MS) recovery(s) of Barium are outside control limits. Spike amount low relative to the sample amount. Refer to lab control or spike blank for recovery information.
- The RPD(s) for the MS and MSD recoveries of Barium are outside control limits for sample MP9011-S2. High RPD due to possible sample matrix or nonhomogeneity.
- The serial dilution RPD(s) for Cadmium, Selenium, Chromium, Lead, Nickel, Zinc are outside control limits for sample MP9011-SD1. Percent difference acceptable due to low initial sample concentration (< 50 times IDL).
- MP9011-SD1 for Chromium: Serial dilution indicates possible matrix interference.
- MP9011-SD1 for Lead: Serial dilution indicates possible matrix interference.
- MP9011-SD1 for Nickel: Serial dilution indicates possible matrix interference.
- MP9011-SD1 for Zinc: Serial dilution indicates possible matrix interference.

## Metals By Method SW846 6020A

**Matrix** SO

**Batch ID:** MP9012

- All samples were digested within the recommended method holding time.
- All samples were analyzed within the recommended method holding time.
- All method blanks for this batch meet method specific criteria.
- Sample(s) D41506-1MS, D41506-1MSD, D41506-1SDL were used as the QC samples for the metals analysis.

## Metals By Method SW846 7471B

**Matrix** SO

**Batch ID:** MP9008

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

## Wet Chemistry By Method ASTM D1498-76M

**Matrix** SO

**Batch ID:** GN17965

- The data for ASTM D1498-76M meets quality control requirements.

## Wet Chemistry By Method SM19 2540B M

**Matrix** SO

**Batch ID:** GN17966

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

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

**Matrix** SO

**Batch ID:** GP8845

- All samples were prepared within the recommended method holding time.
- All samples were analyzed within the recommended method holding time.
- All method blanks for this batch meet method specific criteria.
- Sample(s) D41506-1DUP, D41506-1MS, D41506-1MSD were used as the QC samples for the Chromium, Hexavalent analysis.

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

**Matrix** SO

**Batch ID:** R15381

- The data for SW846 3060A/7196A M meets quality control requirements.
- D41506-1 for Chromium, Trivalent: Calculated as: (Chromium) - (Chromium, Hexavalent)

## Wet Chemistry By Method SW846 9045D

**Matrix** SO

**Batch ID:** GN17963

- The following samples were run outside of holding time for method SW846 9045D: D41506-1

## Wet Chemistry By Method USDA HANDBOOK 60

**Matrix** SO

**Batch ID:** MP9015

- D41506-1A for Sodium Adsorption Ratio: Calculated as:  $(\text{Na meq/L}) / \sqrt{[(\text{Ca meq/L}) + (\text{Mg meq/L})/2]}$

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

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

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

## Summary of Hits

Page 1 of 1

Job Number: D41506  
Account: XTO Energy  
Project: PCU 296-5A  
Collected: 12/03/12

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

### D41506-1 RP SUBLINER(COMP)

TPH-DRO (C10-C28)	33.8	7.4	4.4	mg/kg	SW846-8015B
Arsenic	5.0	0.11		mg/kg	SW846 6020A
Barium	2270	5.5		mg/kg	SW846 6010C
Chromium	23.4	1.1		mg/kg	SW846 6010C
Copper	9.5	1.1		mg/kg	SW846 6010C
Lead	19.9	5.5		mg/kg	SW846 6010C
Nickel	11.9	3.3		mg/kg	SW846 6010C
Zinc	27.7	3.3		mg/kg	SW846 6010C
Specific Conductivity	406	1.0		umhos/cm	SM2510B-1997 MOD
Chromium, Trivalent <sup>a</sup>	23.4	2.1		mg/kg	SW846 3060A/7196A M
Redox Potential Vs H2	168			mv	ASTM D1498-76M
pH	9.93			su	SW846 9045D

### D41506-1A RP SUBLINER(COMP)

Calcium	29.1	2.0		mg/l	SW846 6010C
Magnesium	5.60	1.0		mg/l	SW846 6010C
Sodium	109	2.0		mg/l	SW846 6010C
Sodium Adsorption Ratio <sup>b</sup>	4.84			ratio	USDA HANDBOOK 60

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

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



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## Sample Results

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### Report of Analysis

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**Report of Analysis**

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**Client Sample ID:** RP SUBLINER(COMP)**Lab Sample ID:** D41506-1**Date Sampled:** 12/03/12**Matrix:** SO - Soil**Date Received:** 12/05/12**Method:** SW846 8260B**Percent Solids:** 90.2**Project:** PCU 296-5A

	<b>File ID</b>	<b>DF</b>	<b>Analyzed</b>	<b>By</b>	<b>Prep Date</b>	<b>Prep Batch</b>	<b>Analytical Batch</b>
Run #1	3V22004.D	1	12/07/12	BD	n/a	n/a	V3V1286
Run #2							

	<b>Initial Weight</b>	<b>Final Volume</b>	<b>Methanol Aliquot</b>
Run #1	5.07 g	5.0 ml	100 ul
Run #2			

**Purgeable Aromatics**

<b>CAS No.</b>	<b>Compound</b>	<b>Result</b>	<b>RL</b>	<b>MDL</b>	<b>Units</b>	<b>Q</b>
71-43-2	Benzene	ND	0.060	0.030	mg/kg	
108-88-3	Toluene	ND	0.12	0.060	mg/kg	
100-41-4	Ethylbenzene	ND	0.12	0.023	mg/kg	
1330-20-7	Xylene (total)	ND	0.24	0.12	mg/kg	

<b>CAS No.</b>	<b>Surrogate Recoveries</b>	<b>Run# 1</b>	<b>Run# 2</b>	<b>Limits</b>
2037-26-5	Toluene-D8	97%		64-130%
460-00-4	4-Bromofluorobenzene	87%		62-131%
17060-07-0	1,2-Dichloroethane-D4	115%		70-130%

ND = Not detected MDL - Method Detection Limit

J = Indicates an estimated value

RL = Reporting Limit

B = Indicates analyte found in associated method blank

E = Indicates value exceeds calibration range

N = Indicates presumptive evidence of a compound

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**Report of Analysis**

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<b>Client Sample ID:</b>	RP SUBLINER(COMP)	<b>Date Sampled:</b>	12/03/12
<b>Lab Sample ID:</b>	D41506-1	<b>Date Received:</b>	12/05/12
<b>Matrix:</b>	SO - Soil	<b>Percent Solids:</b>	90.2
<b>Method:</b>	SW846 8270C BY SIM	SW846 3546	
<b>Project:</b>	PCU 296-5A		

	<b>File ID</b>	<b>DF</b>	<b>Analyzed</b>	<b>By</b>	<b>Prep Date</b>	<b>Prep Batch</b>	<b>Analytical Batch</b>
Run #1	3G12516.D	1	12/10/12	DC	12/10/12	OP7075	E3G593
Run #2							

	<b>Initial Weight</b>	<b>Final Volume</b>
Run #1	30.2 g	1.0 ml
Run #2		

**COGCC Table 910-1 PAH List**

<b>CAS No.</b>	<b>Compound</b>	<b>Result</b>	<b>RL</b>	<b>MDL</b>	<b>Units</b>	<b>Q</b>
83-32-9	Acenaphthene	ND	0.0092	0.0048	mg/kg	
120-12-7	Anthracene	ND	0.0092	0.0048	mg/kg	
56-55-3	Benzo(a)anthracene	ND	0.0092	0.0048	mg/kg	
205-99-2	Benzo(b)fluoranthene	ND	0.0092	0.0048	mg/kg	
207-08-9	Benzo(k)fluoranthene	ND	0.0092	0.0048	mg/kg	
50-32-8	Benzo(a)pyrene	ND	0.0092	0.0048	mg/kg	
218-01-9	Chrysene	ND	0.0092	0.0048	mg/kg	
53-70-3	Dibenz(a,h)anthracene	ND	0.0092	0.0048	mg/kg	
206-44-0	Fluoranthene	ND	0.0092	0.0048	mg/kg	
86-73-7	Fluorene	ND	0.0092	0.0048	mg/kg	
193-39-5	Indeno(1,2,3-cd)pyrene	ND	0.0092	0.0048	mg/kg	
91-20-3	Naphthalene	ND	0.013	0.011	mg/kg	
129-00-0	Pyrene	ND	0.0092	0.0048	mg/kg	

<b>CAS No.</b>	<b>Surrogate Recoveries</b>	<b>Run# 1</b>	<b>Run# 2</b>	<b>Limits</b>
4165-60-0	Nitrobenzene-d5	85%		10-159%
321-60-8	2-Fluorobiphenyl	71%		19-131%
1718-51-0	Terphenyl-d14	82%		18-150%

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**

Page 1 of 1

**Client Sample ID:** RP SUBLINER(COMP)**Lab Sample ID:** D41506-1**Date Sampled:** 12/03/12**Matrix:** SO - Soil**Date Received:** 12/05/12**Method:** SW846 8015B**Percent Solids:** 90.2**Project:** PCU 296-5A

	<b>File ID</b>	<b>DF</b>	<b>Analyzed</b>	<b>By</b>	<b>Prep Date</b>	<b>Prep Batch</b>	<b>Analytical Batch</b>
Run #1	GB18764.D	1	12/06/12	SK	n/a	n/a	GGB1023
Run #2							

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

<b>CAS No.</b>	<b>Compound</b>	<b>Result</b>	<b>RL</b>	<b>MDL</b>	<b>Units</b>	<b>Q</b>
	TPH-GRO (C6-C10)	ND	12	6.0	mg/kg	
<b>CAS No.</b>	<b>Surrogate Recoveries</b>	<b>Run# 1</b>	<b>Run# 2</b>	<b>Limits</b>		
120-82-1	1,2,4-Trichlorobenzene	90%		60-140%		

ND = Not detected MDL - Method Detection Limit

J = Indicates an estimated value

RL = Reporting Limit

B = Indicates analyte found in associated method blank

E = Indicates value exceeds calibration range

N = Indicates presumptive evidence of a compound

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**Report of Analysis**

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**Client Sample ID:** RP SUBLINER(COMP)**Lab Sample ID:** D41506-1**Date Sampled:** 12/03/12**Matrix:** SO - Soil**Date Received:** 12/05/12**Method:** SW846-8015B SW846 3546**Percent Solids:** 90.2**Project:** PCU 296-5A

	<b>File ID</b>	<b>DF</b>	<b>Analyzed</b>	<b>By</b>	<b>Prep Date</b>	<b>Prep Batch</b>	<b>Analytical Batch</b>
Run #1	FD20243.D	1	12/07/12	AV	12/06/12	OP7061	GFD1015
Run #2							

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

<b>CAS No.</b>	<b>Compound</b>	<b>Result</b>	<b>RL</b>	<b>MDL</b>	<b>Units</b>	<b>Q</b>
----------------	-----------------	---------------	-----------	------------	--------------	----------

TPH-DRO (C10-C28)	33.8	7.4	4.4	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	68%		35-130%
---------	-------------	-----	--	---------

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

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

**Report of Analysis**

Page 1 of 1

**Client Sample ID:** RP SUBLINER(COMP)**Lab Sample ID:** D41506-1**Matrix:** SO - Soil**Date Sampled:** 12/03/12**Date Received:** 12/05/12**Percent Solids:** 90.2**Project:** PCU 296-5A**Metals Analysis**

Analyte	Result	RL	Units	DF	Prep	Analyzed By	Method	Prep Method
Arsenic	5.0	0.11	mg/kg	5	12/06/12	12/07/12 JM	SW846 6020A <sup>2</sup>	SW846 3050B <sup>7</sup>
Barium	2270	5.5	mg/kg	5	12/06/12	12/10/12 JB	SW846 6010C <sup>4</sup>	SW846 3050B <sup>6</sup>
Cadmium	< 1.1	1.1	mg/kg	1	12/06/12	12/07/12 JB	SW846 6010C <sup>3</sup>	SW846 3050B <sup>6</sup>
Chromium	23.4	1.1	mg/kg	1	12/06/12	12/07/12 JB	SW846 6010C <sup>3</sup>	SW846 3050B <sup>6</sup>
Copper	9.5	1.1	mg/kg	1	12/06/12	12/07/12 JB	SW846 6010C <sup>3</sup>	SW846 3050B <sup>6</sup>
Lead	19.9	5.5	mg/kg	1	12/06/12	12/07/12 JB	SW846 6010C <sup>3</sup>	SW846 3050B <sup>6</sup>
Mercury	< 0.092	0.092	mg/kg	1	12/06/12	12/06/12 JB	SW846 7471B <sup>1</sup>	SW846 7471B <sup>5</sup>
Nickel	11.9	3.3	mg/kg	1	12/06/12	12/07/12 JB	SW846 6010C <sup>3</sup>	SW846 3050B <sup>6</sup>
Selenium	< 5.5	5.5	mg/kg	1	12/06/12	12/07/12 JB	SW846 6010C <sup>3</sup>	SW846 3050B <sup>6</sup>
Silver	< 3.3	3.3	mg/kg	1	12/06/12	12/07/12 JB	SW846 6010C <sup>3</sup>	SW846 3050B <sup>6</sup>
Zinc	27.7	3.3	mg/kg	1	12/06/12	12/07/12 JB	SW846 6010C <sup>3</sup>	SW846 3050B <sup>6</sup>

- (1) Instrument QC Batch: MA3057
- (2) Instrument QC Batch: MA3061
- (3) Instrument QC Batch: MA3062
- (4) Instrument QC Batch: MA3069
- (5) Prep QC Batch: MP9008
- (6) Prep QC Batch: MP9011
- (7) Prep QC Batch: MP9012

RL = Reporting Limit

**Report of Analysis**

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**Client Sample ID:** RP SUBLINER(COMP)**Lab Sample ID:** D41506-1**Matrix:** SO - Soil**Project:** PCU 296-5A**Date Sampled:** 12/03/12**Date Received:** 12/05/12**Percent Solids:** 90.2**General Chemistry**

Analyte	Result	RL	Units	DF	Analyzed	By	Method
<b>prep: DEPT.OF AG, BOOK N9</b>							
Specific Conductivity	406	1.0	umhos/cm	1	12/07/12	JD	SM2510B-1997 MOD
Chromium, Hexavalent	< 1.0	1.0	mg/kg	1	12/07/12	KB	SW846 3060A/7196A
Chromium, Trivalent <sup>a</sup>	23.4	2.1	mg/kg	1	12/07/12 15:49	JB	SW846 3060A/7196A M
Redox Potential Vs H2	168		mv	1	12/05/12	JD	ASTM D1498-76M
Solids, Percent	90.2		%	1	12/06/12	SWT	SM19 2540B M
pH	9.93		su	1	12/05/12 15:00	JD	SW846 9045D

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

RL = Reporting Limit

**Report of Analysis**

Page 1 of 1

4.2  
4**Client Sample ID:** RP SUBLINER(COMP)**Lab Sample ID:** D41506-1A**Matrix:** SO - Soil**Date Sampled:** 12/03/12**Date Received:** 12/05/12**Percent Solids:** 90.2**Project:** PCU 296-5A**SAR Metals Analysis**

Analyte	Result	RL	Units	DF	Prep	Analyzed By	Method	Prep Method
Calcium	29.1	2.0	mg/l	1	12/06/12	12/10/12 JB	SW846 6010C <sup>1</sup>	SW846 3010A/M <sup>2</sup>
Magnesium	5.60	1.0	mg/l	1	12/06/12	12/10/12 JB	SW846 6010C <sup>1</sup>	SW846 3010A/M <sup>2</sup>
Sodium	109	2.0	mg/l	1	12/06/12	12/10/12 JB	SW846 6010C <sup>1</sup>	SW846 3010A/M <sup>2</sup>

(1) Instrument QC Batch: MA3069

(2) Prep QC Batch: MP9015

RL = Reporting Limit

**Report of Analysis**

Page 1 of 1

**Client Sample ID:** RP SUBLINER(COMP)**Lab Sample ID:** D41506-1A**Matrix:** SO - Soil**Project:** PCU 296-5A**Date Sampled:** 12/03/12**Date Received:** 12/05/12**Percent Solids:** 90.2**General Chemistry**

Analyte	Result	RL	Units	DF	Analyzed	By	Method
Sodium Adsorption Ratio <sup>a</sup>	4.84		ratio	1	12/10/12 10:13	JB	USDA HANDBOOK 60

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

RL = Reporting Limit



## Misc. Forms

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5

### Custody Documents and Other Forms

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Includes the following where applicable:

- Chain of Custody



## CHAIN OF CUSTODY

PAGE 1 OF 1

4036 Youngfield Street, Wheat Ridge, CO 80033  
TEL. 303-425-5021 FAX. 303-425-6854  
www.accutest.com

FED-EX Tracking #	Bottle Order Control #
	D41506

Client / Reporting Information		Project Information		Requested Analysis (see TEST CODE sheet)										Matrix Codes						
Company Name <b>KRW Consulting</b>	Project Name: <b>XTO PCW 296-5A</b>	Street Address <b>8000 West 14th Street, Suite 200</b>	Street	Billing Information ( If different from Report to)																
City <b>Lakewood, CO 80214</b>	City	State	Company Name <b>XTO Energy</b>																	
Project Contact <b>Dwayne Knudson</b>	Project # <b>1210-04</b>	Client Purchase Order # <b>21458 CR 5</b>	Street Address <b>Rifle, CO 81650</b>																	
Phone # <b>970-488-1098</b>	City	State	Attention <b>Jessica Dooling</b>																	
Sampler(s) Name(s) <b>DAVID SANDERS</b>	Project Manager <b>Joe Hess</b>	Phone # <b>970-488-1098</b>	Attention <b>Jessica Dooling</b>																	
Collection		Date	Time	Sampled by	Matrix	# of bottles	HCl	NaOH	HNO3	H2SO4	NONE	Li/Water	NaOH	ENCORE	Bottles	Number of preserved Bottles				
Accutest Sample #	Field ID / Point of Collection <b>RP SUBLINER (COMP)</b>	MEOH/DI Vial # <b>12-3-12</b>	<b>10:20</b>	<b>DS</b>	<b>SO</b>	<b>5</b>					X					<b>X</b>				
T-910																				
J20104																				
Turnaround Time ( Business days)		Data Deliverable Information										Comments / Special Instructions								
<input type="checkbox"/> Std. 10 Business Days <input checked="" type="checkbox"/> Std. 5 Business Days (By contract only) <input type="checkbox"/> 3 Day Emergency <input type="checkbox"/> 2 Day Emergency <input type="checkbox"/> 1 Day Emergency		Approved By (Accutest PM): / Date: <hr/> <hr/> <hr/> <hr/> <hr/>										<input type="checkbox"/> Commercial "A" (Level 1) <input type="checkbox"/> Commercial "B" (Level 2) <input type="checkbox"/> COMMNB <input type="checkbox"/> COMMNB+						<input type="checkbox"/> State Forms Required <input type="checkbox"/> Send Forms to State <input type="checkbox"/> Report by Fax <input checked="" type="checkbox"/> Report by PDF ONLY <input type="checkbox"/> EDD Format	Please email to: <b>KRW Piceance Team</b>	
Emergency & Rush T/A data available VIA LabLink		Commercial "A" = Results Only Commercial "B" = Results + QC Summary Commercial BN = Results/QC/Narrative (+ = chromatograms)																		
Sample Custody must be documented below each time samples change possession, including courier delivery.																				
Relinquished by Sampler: <b>1 Lori Albinson</b>	Date/Time: <b>10/14/12 10:30</b>	Received By: <b>1</b>	Relinquished By: <b>2</b>	Date/Time: <b>10/14/12 10:30</b>	Received By: <b>2</b>															
Relinquished by Sampler: <b>3</b>	Date/Time: <b>10/15/12 10:00</b>	Received By: <b>3</b>	Relinquished By: <b>4</b>	Date/Time: <b>10/15/12 10:00</b>	Received By: <b>4</b>															
Relinquished by: <b>5</b>	Date/Time: <b>10/15/12 10:00</b>	Received By: <b>5</b>	Custody Seal # <b>5</b>	<input type="checkbox"/> Intact <input type="checkbox"/> Not Intact	Preserved where applicable <b>5</b>	On Ice <b>5</b>	Cooler Temp. <b>32°</b>													

D41506: Chain of Custody

Page 1 of 2



## Accutest Laboratories Sample Receipt Summary

Accutest Job Number: D41506

Client: KRW

Immediate Client Services Action Required: No

Date / Time Received: 12/5/2012 10:00:00 AM

No. Coolers:

1

Client Service Action Required at Login: No

Project: XTO PCU 296-5A

Airbill #'s:

**Cooler Security**Y or N

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

**Cooler Temperature**Y or N

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

**Quality Control Preservation**Y or N

N/A

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

**Sample Integrity - Documentation**Y or N

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

**Sample Integrity - Condition**Y or N

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

**Sample Integrity - Instructions**Y or N

N/A

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

Comments

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

5.1

5

**D41506: Chain of Custody****Page 2 of 2**



## GC/MS Volatiles

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### QC Data Summaries

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Includes the following where applicable:

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



## Method Blank Summary

Page 1 of 1

**Job Number:** D41506  
**Account:** XTOKRWR XTO Energy  
**Project:** PCU 296-5A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
V3V1286-MB	3V21999.D	1	12/06/12	BD	n/a	n/a	V3V1286

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

**Method:** SW846 8260B

D41506-1

6.11

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

CAS No.	Surrogate Recoveries	Limits
2037-26-5	Toluene-D8	101%
460-00-4	4-Bromofluorobenzene	84%
17060-07-0	1,2-Dichloroethane-D4	124%

**Blank Spike Summary**

**Job Number:** D41506  
**Account:** XTOKWR XTO Energy  
**Project:** PCU 296-5A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
V3V1286-BS	3V22000.D	1	12/06/12	BD	n/a	n/a	V3V1286

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

D41506-1

CAS No.	Compound	Spike ug/kg	BSP ug/kg	BSP %	Limits
71-43-2	Benzene	50	49.9	100	70-130
100-41-4	Ethylbenzene	50	47.3	95	70-130
108-88-3	Toluene	50	45.4	91	70-130
1330-20-7	Xylene (total)	150	142	95	70-130

CAS No.	Surrogate Recoveries	BSP	Limits
2037-26-5	Toluene-D8	100%	64-130%
460-00-4	4-Bromofluorobenzene	98%	62-131%
17060-07-0	1,2-Dichloroethane-D4	125%	70-130%

\* = Outside of Control Limits.

# Matrix Spike/Matrix Spike Duplicate Summary

Page 1 of 1

Job Number: D41506

Account: XTOKWR XTO Energy

Project: PCU 296-5A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
D41506-1MS	3V22002.D	1	12/07/12	BD	n/a	n/a	V3V1286
D41506-1MSD	3V22003.D	1	12/07/12	BD	n/a	n/a	V3V1286
D41506-1	3V22004.D	1	12/07/12	BD	n/a	n/a	V3V1286

The QC reported here applies to the following samples:

Method: SW846 8260B

D41506-1

CAS No.	Compound	D41506-1		Spike	MS	MS	MSD	MSD	Limits	
		ug/kg	Q	ug/kg	ug/kg	%	ug/kg	%	RPD	Rec/RPD
71-43-2	Benzene	ND		3000	3130	104	3030	101	3	64-139/30
100-41-4	Ethylbenzene	ND		3000	3010	100	2900	97	4	68-136/30
108-88-3	Toluene	ND		3000	2720	91	2650	88	3	60-130/30
1330-20-7	Xylene (total)	ND		9010	9080	101	8830	98	3	58-142/30

CAS No.	Surrogate Recoveries	MS	MSD	D41506-1	Limits
2037-26-5	Toluene-D8	97%	95%	97%	64-130%
460-00-4	4-Bromofluorobenzene	97%	98%	87%	62-131%
17060-07-0	1,2-Dichloroethane-D4	122%	114%	115%	70-130%

\* = Outside of Control Limits.

6.3.1  
6



GC/MS Volatiles

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Raw Data

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7

## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\V3120612.S\  
 Data File : 3V22004.D  
 Acq On : 7 Dec 2012 1:43 am  
 Operator : BRETD  
 Sample : D41506-1  
 Misc : MS5066,V3V1286,5.072,,100,5,1  
 ALS Vial : 31 Sample Multiplier: 1

Quant Time: Dec 07 11:53:31 2012  
 Quant Method : C:\msdchem\1\METHODS\V3AP1277TVH1277SOIL.M  
 Quant Title : 8260  
 QLast Update : Wed Nov 28 14:20:19 2012  
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
2) Pentafluorobenzene	11.863	168	124817	50.00	ug/l	0.00
35) 1,4-Difluorobenzene	12.655	114	221305	50.00	ug/l	0.00
53) Chlorobenzene-d5	15.292	117	255859	50.00	ug/l	0.00
74) 1,4-Dichlorobenzene-d4	17.284	152	133689	50.00	ug/l	0.00

System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	12.247	102	17367	57.32	ug/l	0.00
Spiked Amount 50.000	Range 70 - 130		Recovery	=	114.64%	
61) Toluene-d8	14.050	98	298056	48.32	ug/l	0.00
Spiked Amount 50.000	Range 70 - 130		Recovery	=	96.64%	
69) 4-Bromofluorobenzene	16.244	95	110817	43.71	ug/l	0.00
Spiked Amount 50.000	Range 70 - 130		Recovery	=	87.42%	

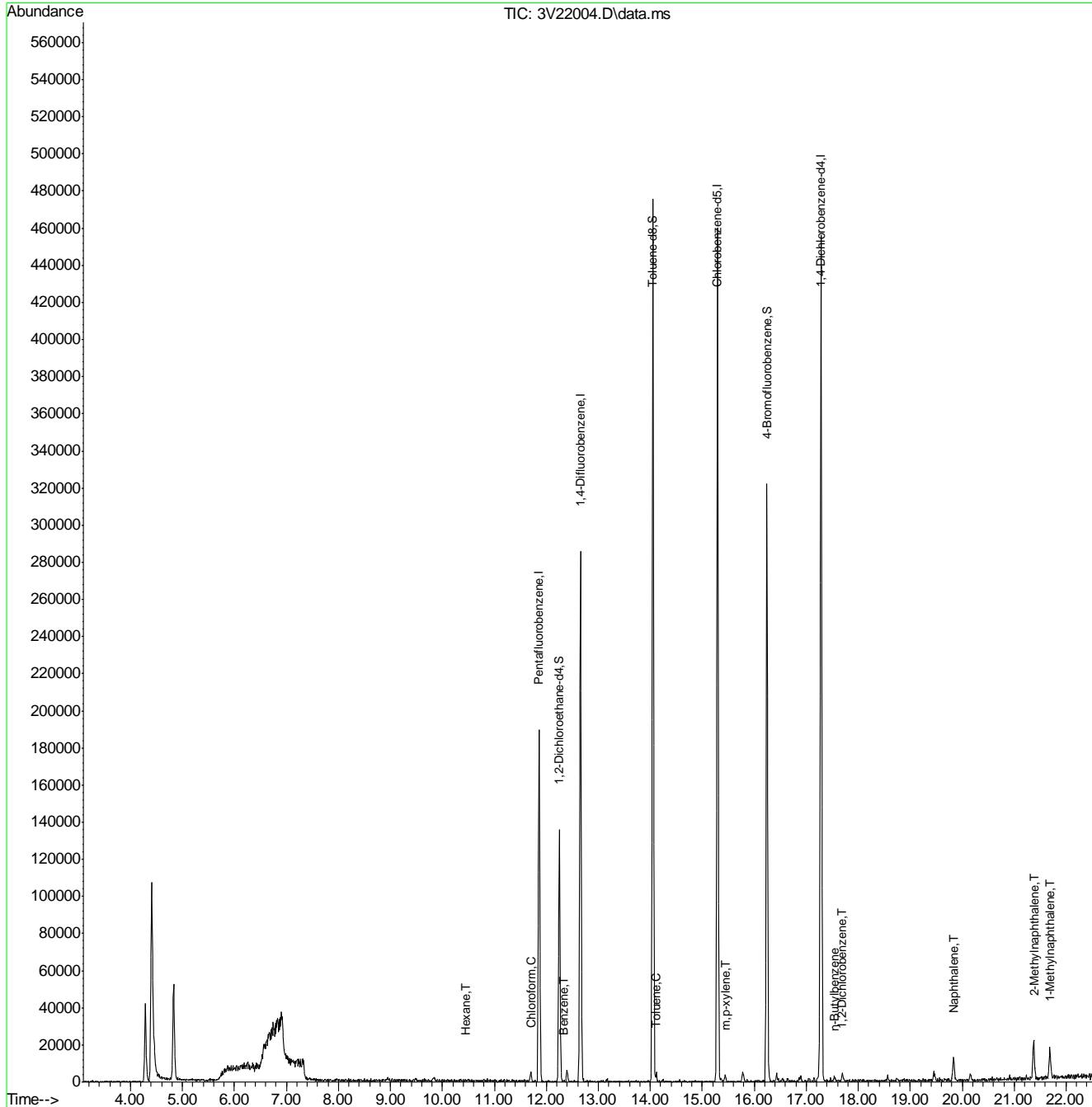
Target Compounds					Qvalue
1) TVH-Gasoline	13.285	TIC	-9125m	158.76	ug/l
29) Chloroform	11.702	83	4229	1.33	ug/l 91
41) Hexane	10.458	57	132	0.04	ug/l 100
50) Benzene	12.340	78	196	0.03	ug/l 100
62) Toluene	14.114	92	1818	0.33	ug/l 100
72) m,p-xylene	15.446	106	1352	0.32	ug/l 99
87) 1,2-Dichlorobenzene	17.688	146	1906	0.43	ug/l # 93
88) n-Butylbenzene	17.543	91	2407	0.31	ug/l # 85
91) Naphthalene	19.843	128	15427	5.25	ug/l 100
94) 2-Methylnaphthalene	21.377	142	15460	5.94	ug/l 98
95) 1-Methylnaphthalene	21.688	142	12595	5.27	ug/l 98

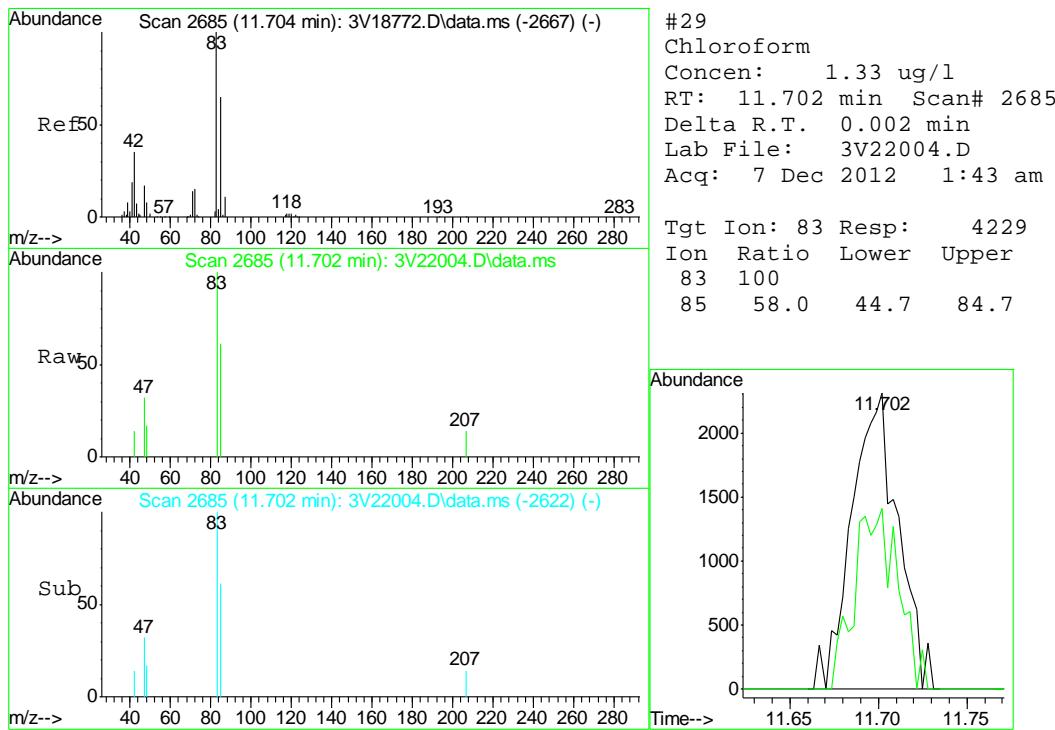
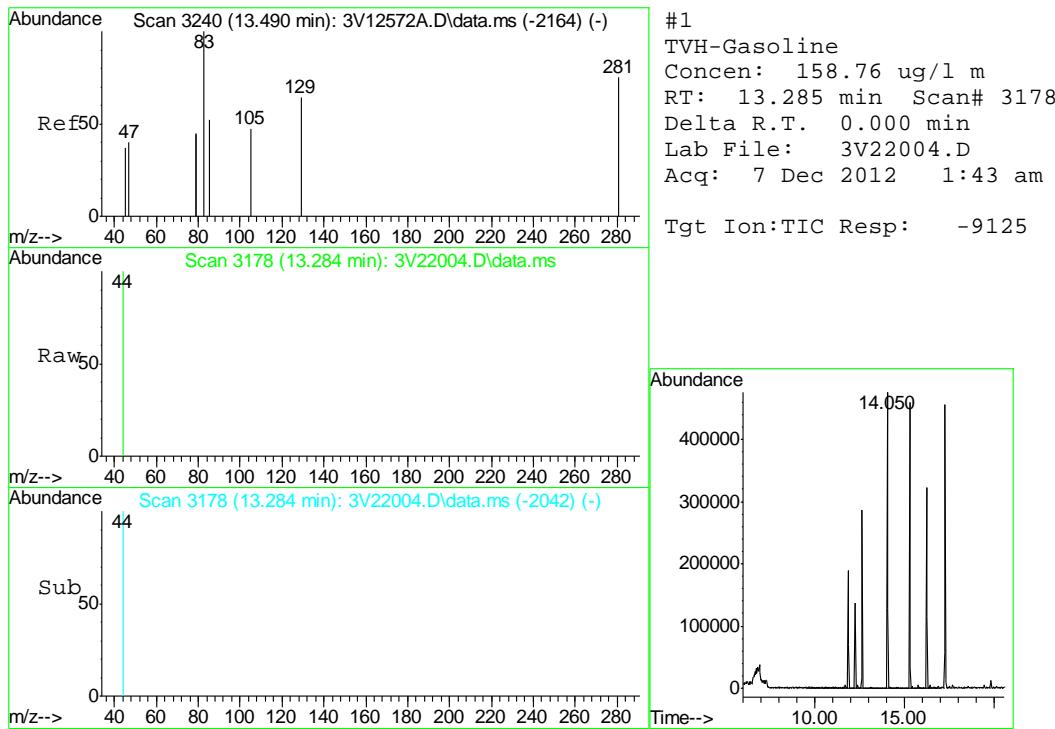
(#) = qualifier out of range (m) = manual integration (+) = signals summed

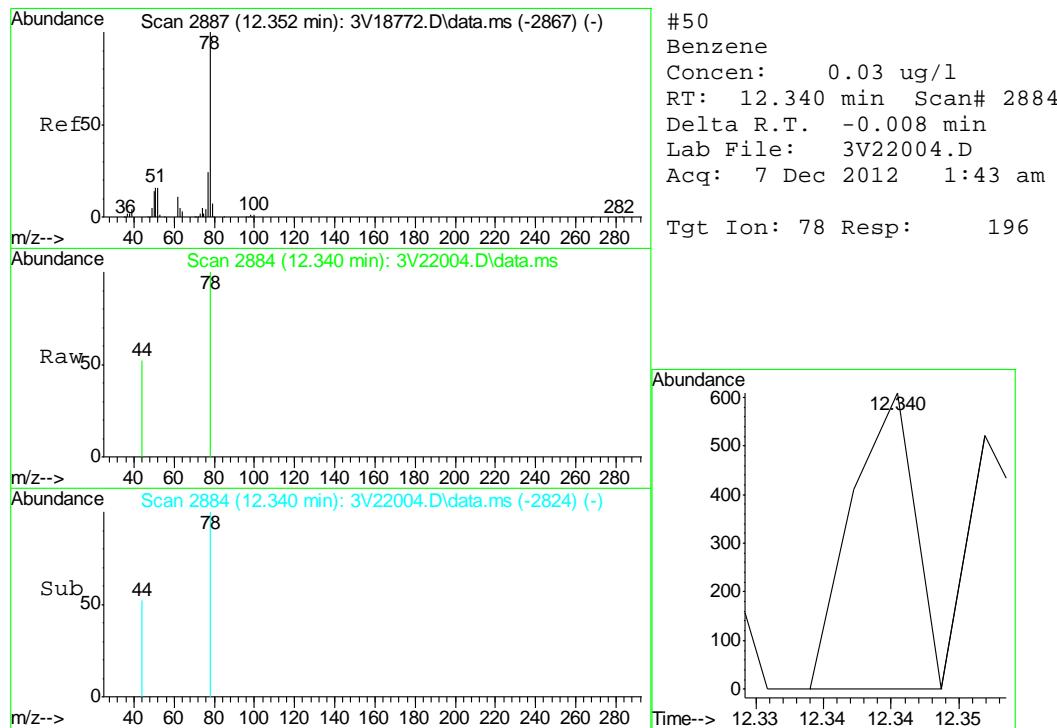
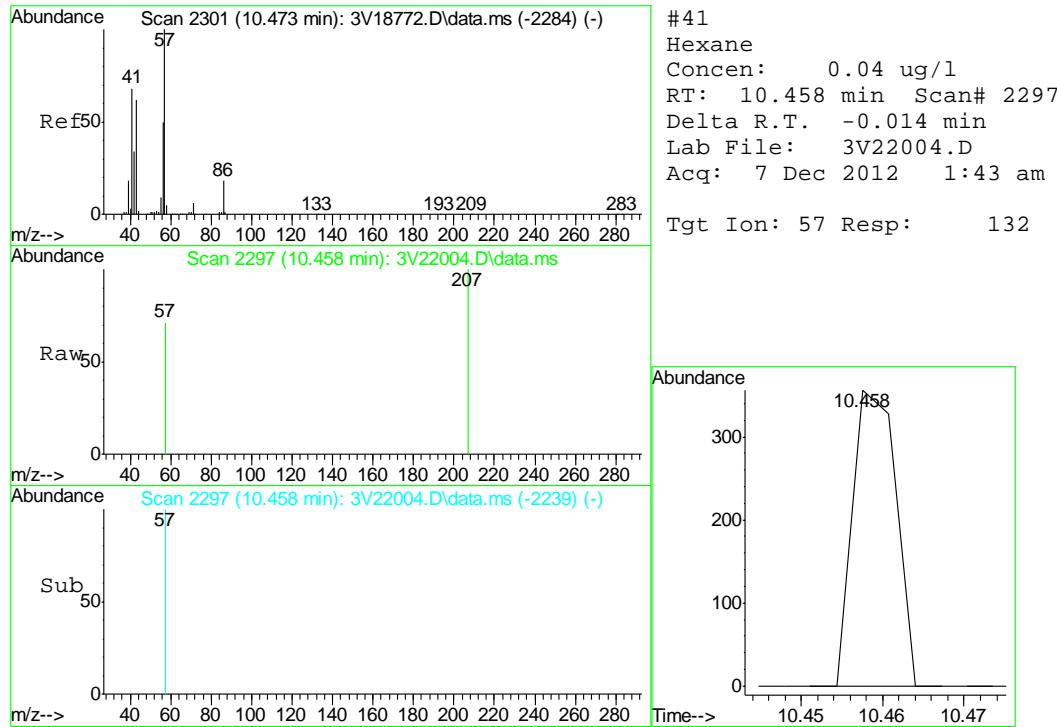
## Quantitation Report (QT Reviewed)

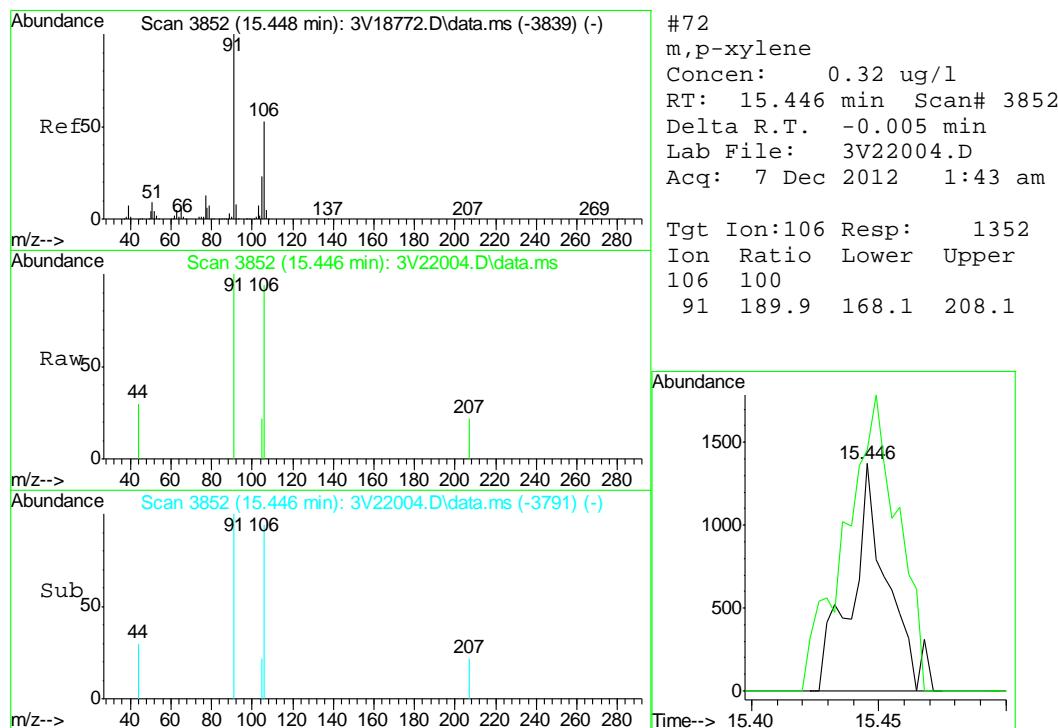
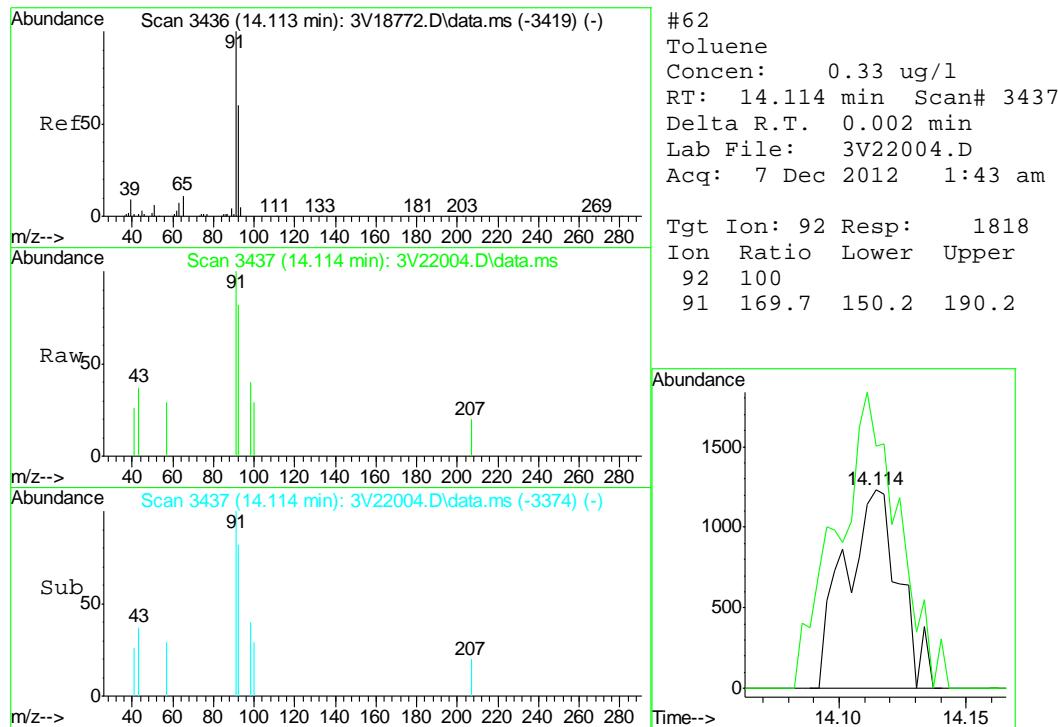
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 Operator : BRETD  
 Sample : D41506-1  
 Misc : MS5066,V3V1286,5.072,,100,5,1  
 ALS Vial : 31 Sample Multiplier: 1

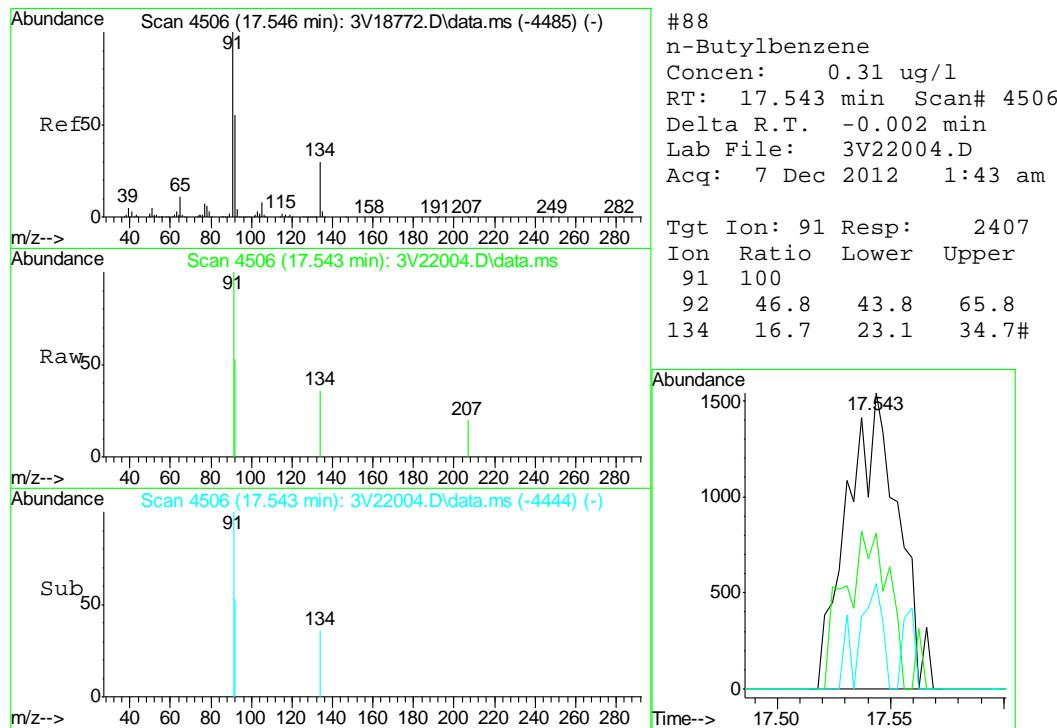
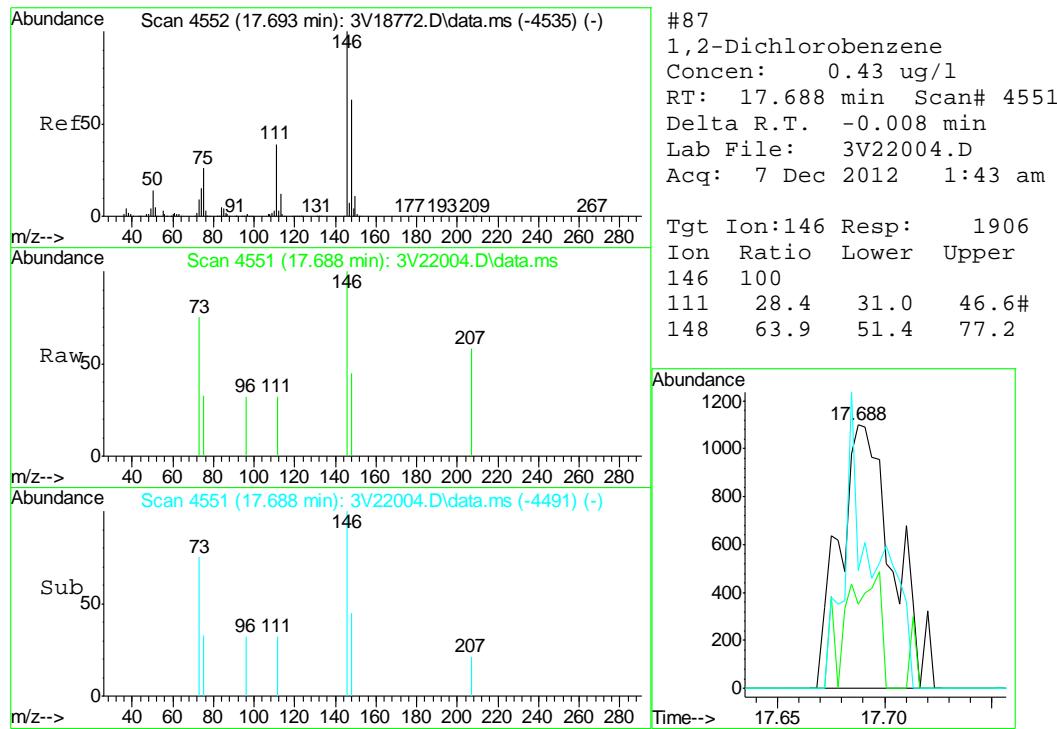
Quant Time: Dec 07 11:53:31 2012  
 Quant Method : C:\msdchem\1\METHODS\V3AP1277TVH1277SOIL.M  
 Quant Title : 8260  
 QLast Update : Wed Nov 28 14:20:19 2012  
 Response via : Initial Calibration

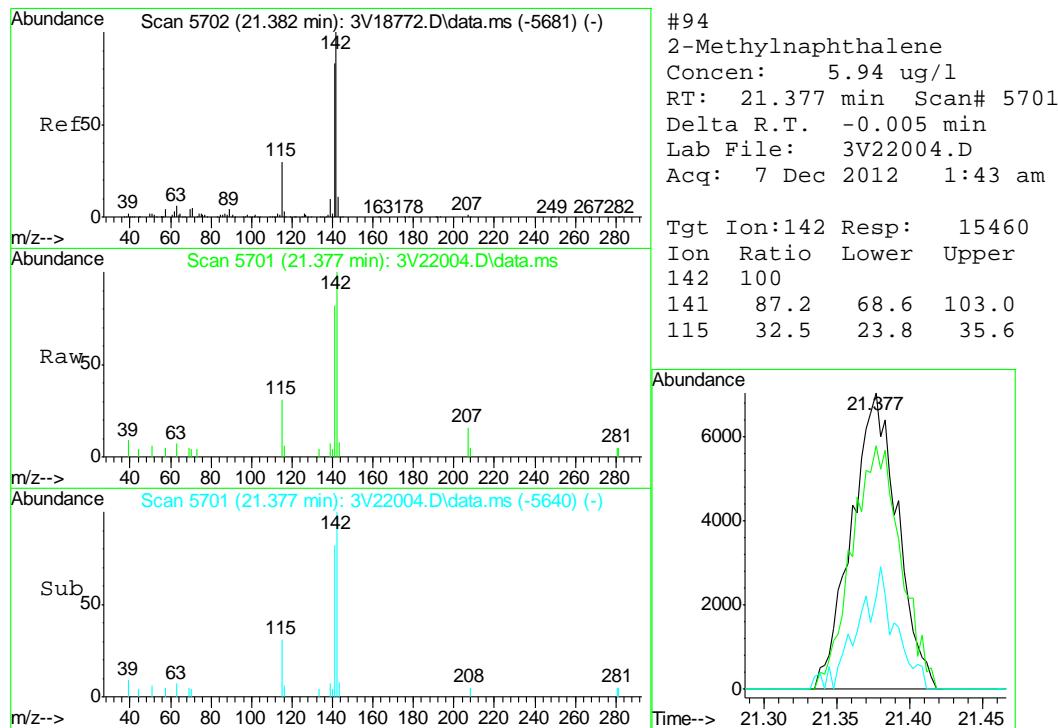
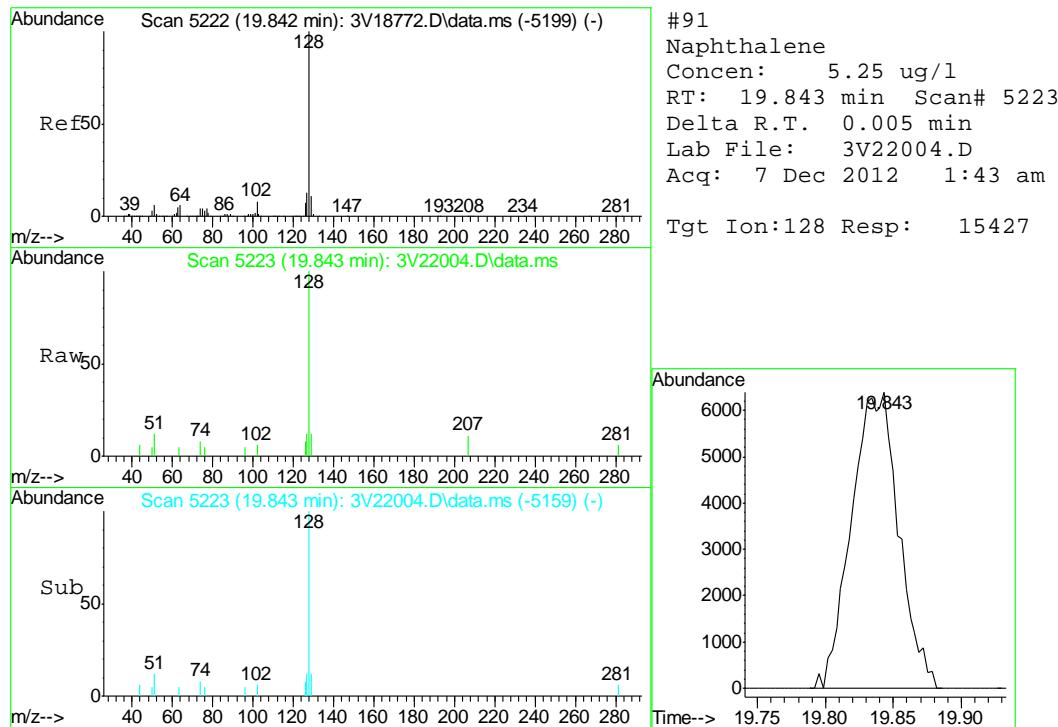


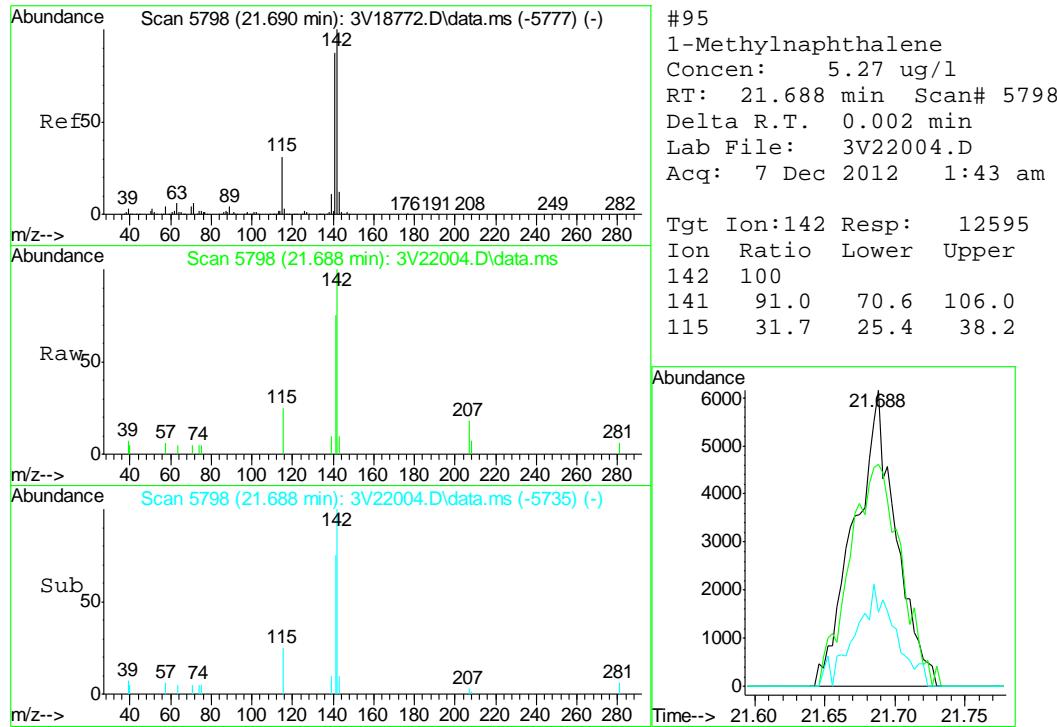












## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\V3120612.S\  
 Data File : 3V21999.D  
 Acq On : 6 Dec 2012 11:07 pm  
 Operator : BRETD  
 Sample : MB  
 Misc : MS5066,V3V1286,5.00,,100,5,1  
 ALS Vial : 26 Sample Multiplier: 1

Quant Time: Dec 07 11:33:59 2012  
 Quant Method : C:\msdchem\1\METHODS\V3AP1277TVH1277SOIL.M  
 Quant Title : 8260  
 QLast Update : Wed Nov 28 14:20:19 2012  
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
2) Pentafluorobenzene	11.864	168	108264	50.00	ug/l	0.00
35) 1,4-Difluorobenzene	12.660	114	197587	50.00	ug/l	0.00
53) Chlorobenzene-d5	15.293	117	214080	50.00	ug/l	0.00
74) 1,4-Dichlorobenzene-d4	17.282	152	112638	50.00	ug/l	0.00

System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	12.249	102	16296	62.00	ug/l	0.00
Spiked Amount 50.000	Range 70 - 130		Recovery	=	124.00%	
61) Toluene-d8	14.049	98	261414	50.65	ug/l	0.00
Spiked Amount 50.000	Range 70 - 130		Recovery	=	101.30%	
69) 4-Bromofluorobenzene	16.240	95	89291	42.10	ug/l	0.00
Spiked Amount 50.000	Range 70 - 130		Recovery	=	84.20%	

Target Compounds					Qvalue
1) TVH-Gasoline	13.285	TIC	60208m	161.01	ug/l
41) Hexane	10.469	57	148	0.05	ug/l 100
50) Benzene	12.352	78	73	0.01	ug/l 100
62) Toluene	14.110	92	1066	0.23	ug/l # 83
71) Styrene	15.781	104	191	0.18	ug/l 81
91) Naphthalene	19.839	128	7801	4.47	ug/l 100
94) 2-Methylnaphthalene	21.378	142	7469	3.40	ug/l 98
95) 1-Methylnaphthalene	21.693	142	5651	2.81	ug/l # 88

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

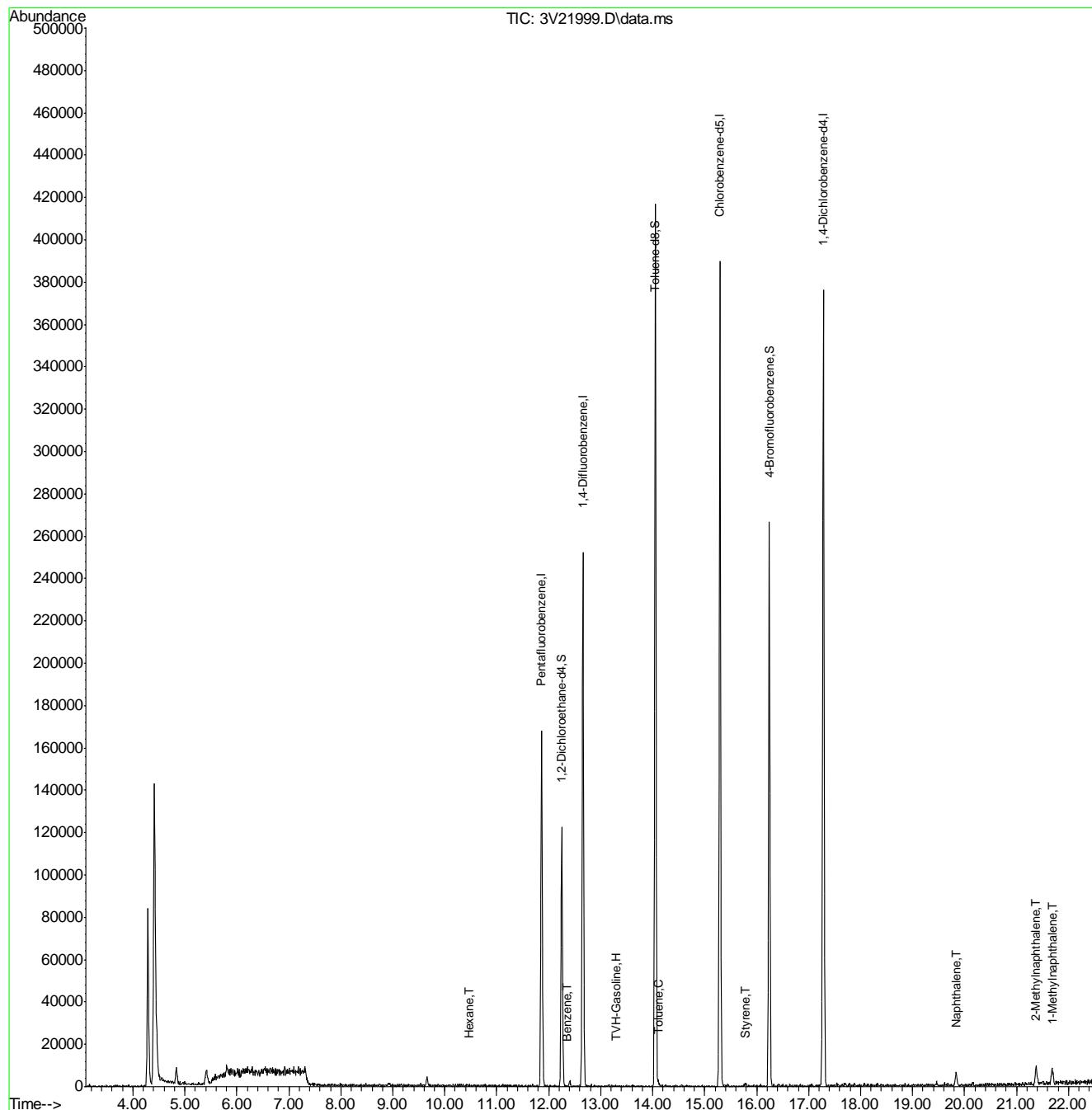
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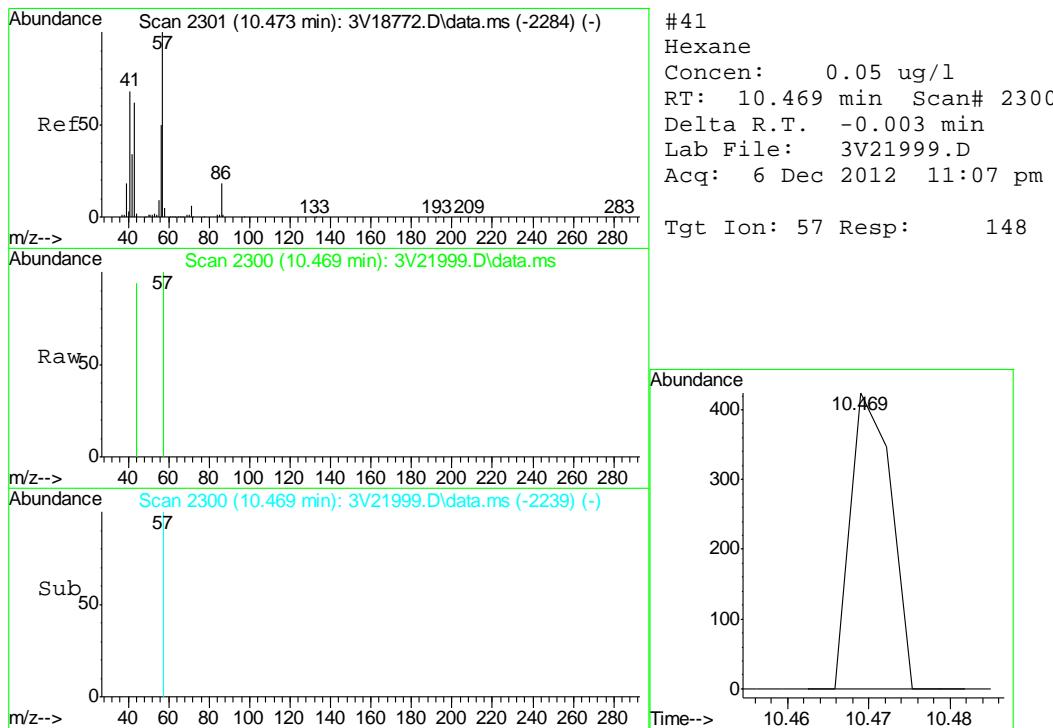
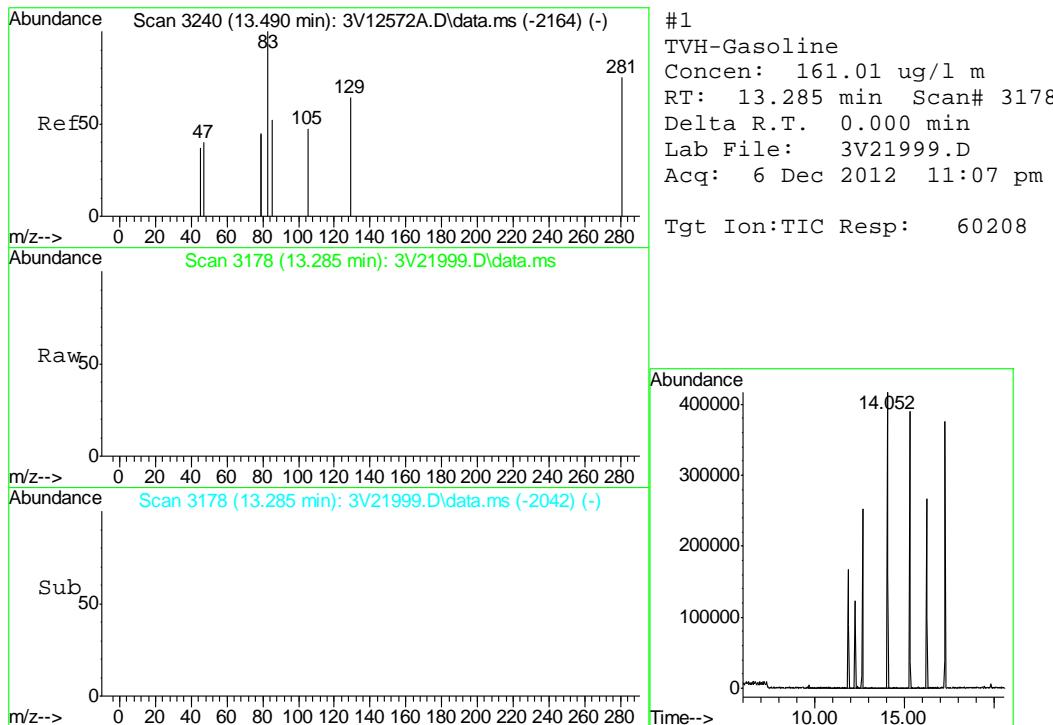
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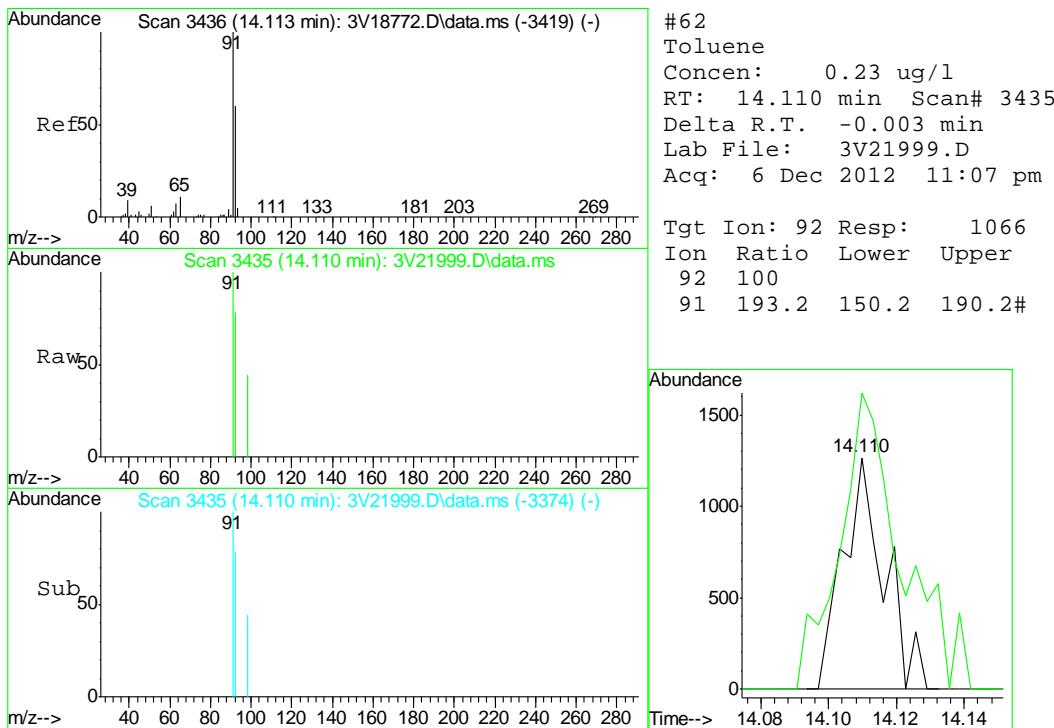
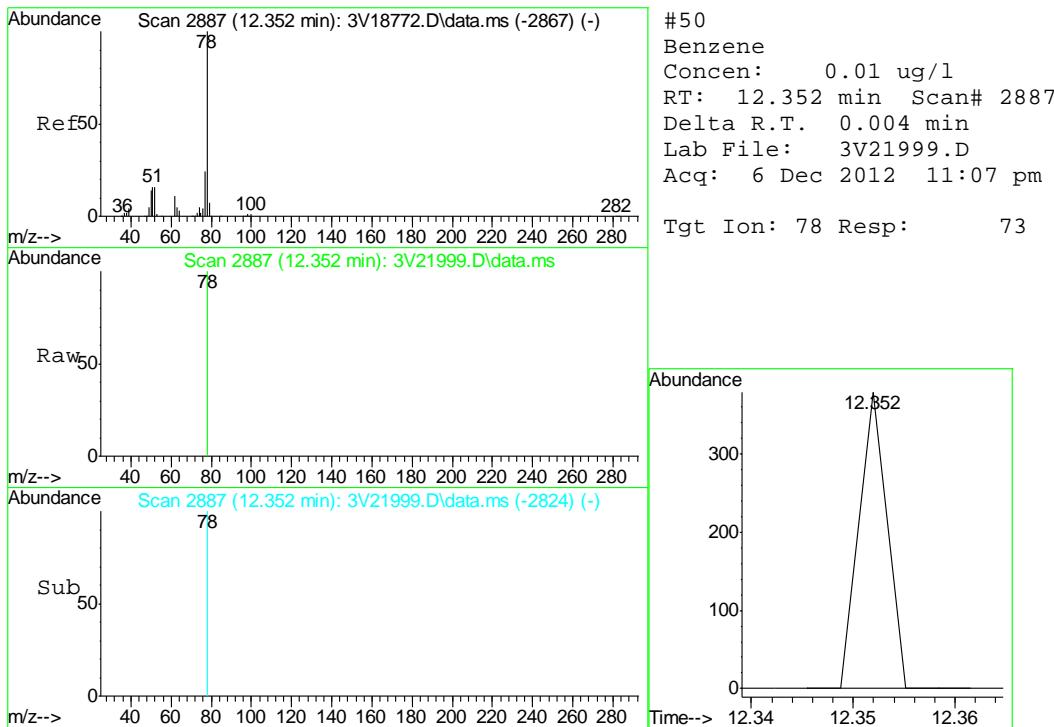
## Quantitation Report (QT Reviewed)

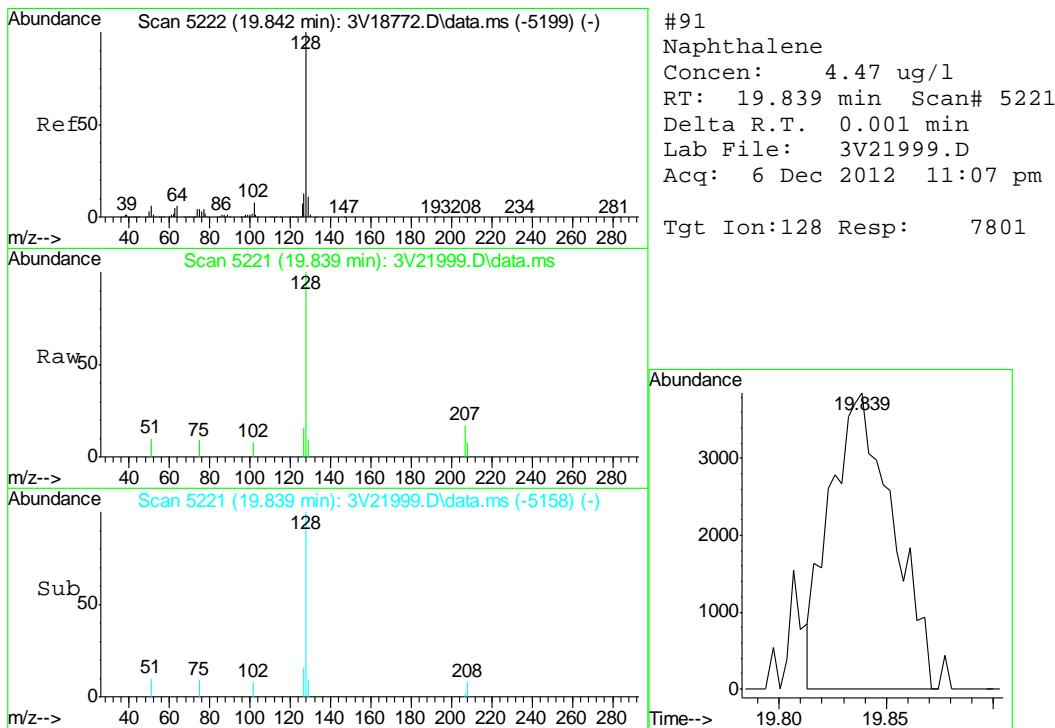
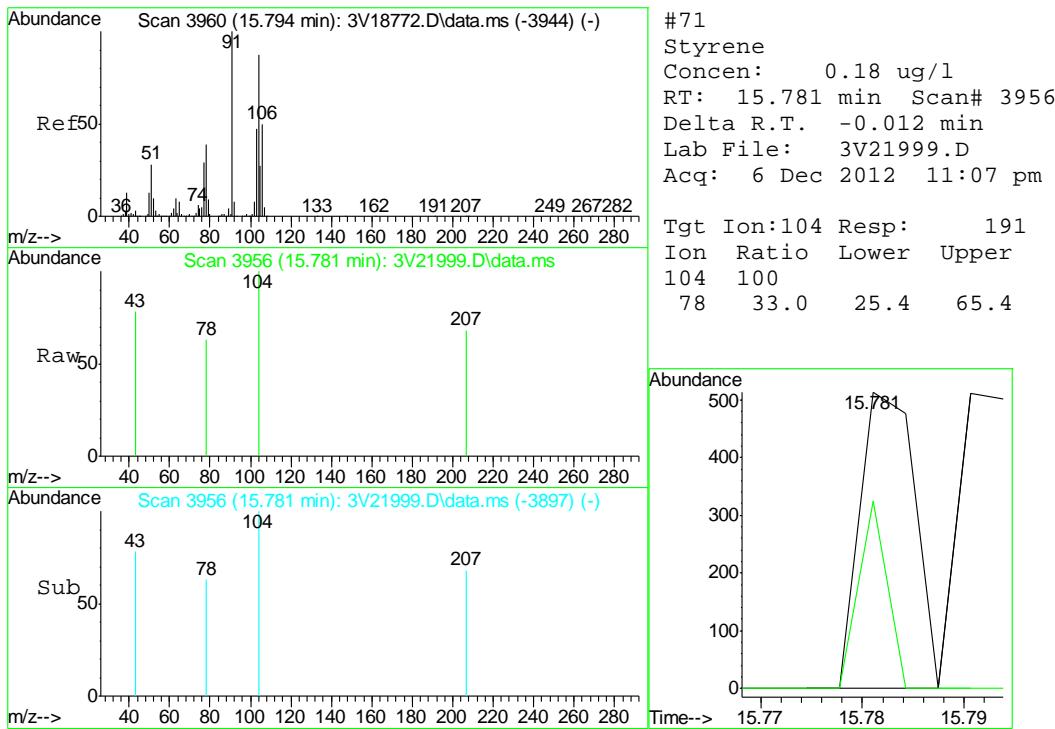
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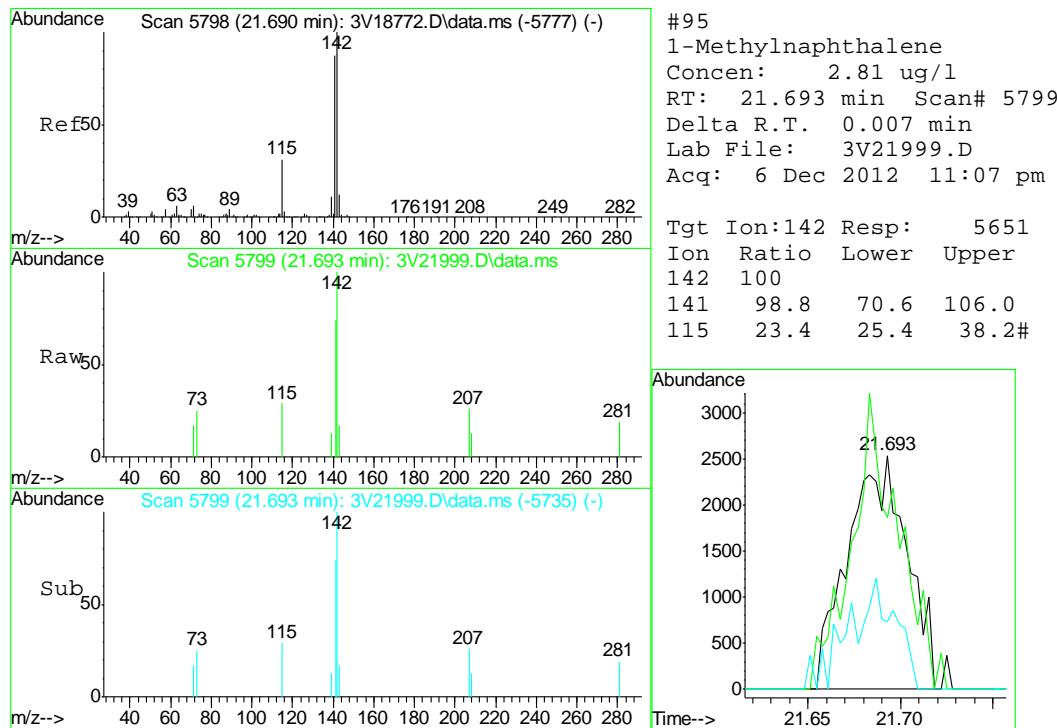
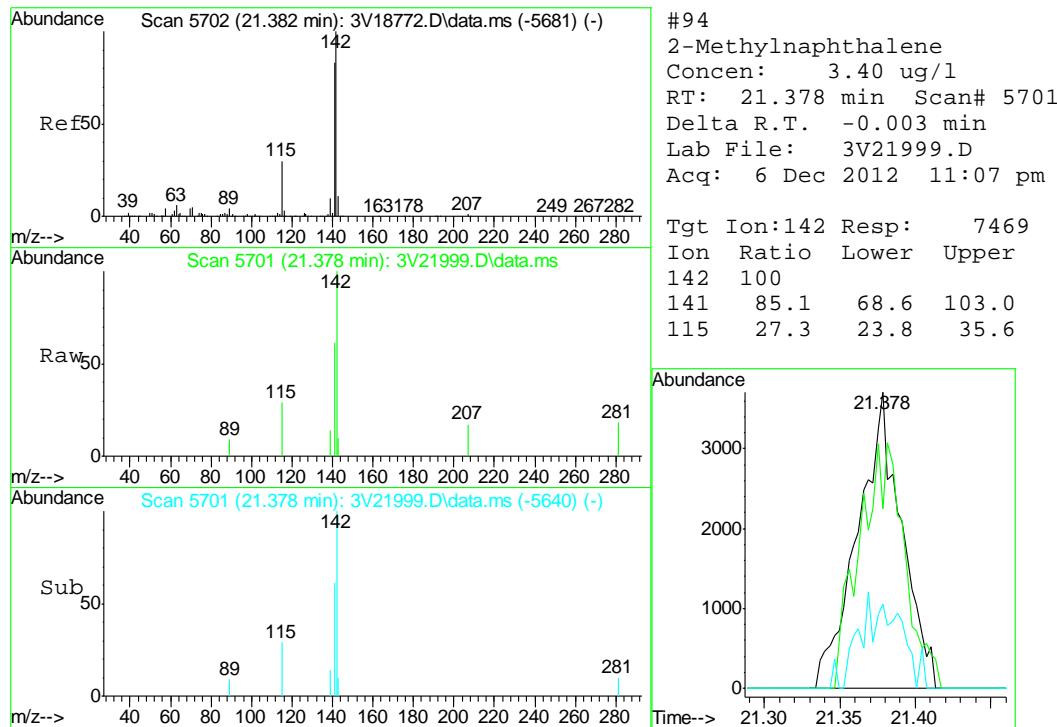
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 Quant Method : C:\msdchem\1\METHODS\V3AP1277TVH1277SOIL.M  
 Quant Title : 8260  
 QLast Update : Wed Nov 28 14:20:19 2012  
 Response via : Initial Calibration













## GC/MS Semi-volatiles

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### QC Data Summaries

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Includes the following where applicable:

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

**Method Blank Summary**

**Job Number:** D41506  
**Account:** XTOKWR XTO Energy  
**Project:** PCU 296-5A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
OP7075-MB	3G12508.D	1	12/10/12	DC	12/10/12	OP7075	E3G593

The QC reported here applies to the following samples:

**Method:** SW846 8270C BY SIM

D41506-1

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

CAS No.	Surrogate Recoveries	Limits
4165-60-0	Nitrobenzene-d5	88%
321-60-8	2-Fluorobiphenyl	79%
1718-51-0	Terphenyl-d14	91%

## Blank Spike Summary

Page 1 of 1

Job Number: D41506

Account: XTOKWR XTO Energy

Project: PCU 296-5A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
OP7075-BS	3G12509.D	1	12/10/12	DC	12/10/12	OP7075	E3G593

The QC reported here applies to the following samples:

Method: SW846 8270C BY SIM

D41506-1

CAS No.	Compound	Spike ug/kg	BSP ug/kg	BSP %	Limits
83-32-9	Acenaphthene	83.3	74.0	89	68-130
120-12-7	Anthracene	83.3	63.0	76	67-130
56-55-3	Benzo(a)anthracene	83.3	72.7	87	65-130
205-99-2	Benzo(b)fluoranthene	83.3	83.2	100	44-130
207-08-9	Benzo(k)fluoranthene	83.3	66.7	80	56-131
50-32-8	Benzo(a)pyrene	83.3	74.9	90	62-130
218-01-9	Chrysene	83.3	74.6	90	70-130
53-70-3	Dibenzo(a,h)anthracene	83.3	72.2	87	55-130
206-44-0	Fluoranthene	83.3	63.1	76	70-130
86-73-7	Fluorene	83.3	71.5	86	70-130
193-39-5	Indeno(1,2,3-cd)pyrene	83.3	72.0	86	56-130
91-20-3	Naphthalene	83.3	78.0	94	70-130
129-00-0	Pyrene	83.3	76.0	91	70-130

CAS No.	Surrogate Recoveries	BSP	Limits
4165-60-0	Nitrobenzene-d5	97%	10-159%
321-60-8	2-Fluorobiphenyl	76%	19-131%
1718-51-0	Terphenyl-d14	87%	18-150%

\* = Outside of Control Limits.

# Matrix Spike/Matrix Spike Duplicate Summary

Page 1 of 1

Job Number: D41506

Account: XTOKWR XTO Energy

Project: PCU 296-5A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
OP7075-MS	3G12511.D	1	12/10/12	DC	12/10/12	OP7075	E3G593
OP7075-MSD	3G12512.D	1	12/10/12	DC	12/10/12	OP7075	E3G593
D41381-1	3G12510.D	1	12/10/12	DC	12/10/12	OP7075	E3G593

The QC reported here applies to the following samples:

Method: SW846 8270C BY SIM

D41506-1

CAS No.	Compound	D41381-1		Spike	MS	MS	MSD	MSD	RPD	Limits Rec/RPD
		ug/kg	Q	ug/kg	ug/kg	%	ug/kg	%		
83-32-9	Acenaphthene	ND		93.9	77.2	82	74.1	79	4	25-151/30
120-12-7	Anthracene	ND		93.9	71.0	76	69.6	74	2	39-159/30
56-55-3	Benzo(a)anthracene	ND		93.9	80.7	86	79.9	85	1	39-168/30
205-99-2	Benzo(b)fluoranthene	ND		93.9	85.0	90	85.9	92	1	24-163/30
207-08-9	Benzo(k)fluoranthene	ND		93.9	77.1	82	76.7	82	1	10-188/30
50-32-8	Benzo(a)pyrene	ND		93.9	83.5	89	81.1	86	3	32-144/30
218-01-9	Chrysene	ND		93.9	80.3	85	81.0	86	1	43-150/30
53-70-3	Dibenzo(a,h)anthracene	ND		93.9	81.0	86	77.6	83	4	21-152/30
206-44-0	Fluoranthene	ND		93.9	71.8	76	69.7	74	3	36-157/30
86-73-7	Fluorene	ND		93.9	80.0	85	74.9	80	7	10-182/30
193-39-5	Indeno(1,2,3-cd)pyrene	ND		93.9	79.8	85	77.7	83	3	20-154/30
91-20-3	Naphthalene	ND		93.9	77.2	82	73.0	78	6	10-163/30
129-00-0	Pyrene	ND		93.9	83.7	89	83.2	89	1	25-180/30

CAS No.	Surrogate Recoveries	MS	MSD	D41381-1	Limits
4165-60-0	Nitrobenzene-d5	83%	78%	70%	10-159%
321-60-8	2-Fluorobiphenyl	69%	64%	57%	19-131%
1718-51-0	Terphenyl-d14	78%	78%	72%	18-150%

\* = Outside of Control Limits.

8.3.1  
8



## GC/MS Semi-volatiles

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Raw Data

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## Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\121012\  
 Data File : 3g12516.D  
 Acq On : 10 Dec 2012 3:02 pm  
 Operator : DONC  
 Sample : D41506-1  
 Misc : OP7075,E3G593,30.16,,,1,1  
 ALS Vial : 12 Sample Multiplier: 1

Quant Time: Dec 11 13:04:13 2012  
 Quant Method : C:\msdchem\1\METHODS\SIMPE3G586.M  
 Quant Title : PAHSIM BASE  
 QLast Update : Tue Dec 04 08:50:28 2012  
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Naphthalene-d8	5.670	136	152027	4.0000	ug/mL	-0.01
6) Acenaphthene-d10	7.384	164	97440	4.0000	ug/mL	0.00
15) Phenanthrene-d10	8.859	188	163716	4.0000	ug/mL	-0.02
19) Chrysene-d12	11.496	240	113350	4.0000	ug/mL	-0.02
24) Perylene-d12	12.873	264	88802	4.0000	ug/mL	-0.02

System Monitoring Compounds						
2) Nitrobenzene-d5	4.985	82	644794	42.4294	ug/mL	-0.01
Spiked Amount	50.000	Range	25 - 135	Recovery	=	84.86%
7) 2-Fluorobiphenyl	6.723	172	1525264	35.4196	ug/mL	0.00
Spiked Amount	50.000	Range	25 - 135	Recovery	=	70.84%
21) Terphenyl-d14	10.450	244	684818	41.0446	ug/mL	-0.02
Spiked Amount	50.000	Range	25 - 135	Recovery	=	82.08%

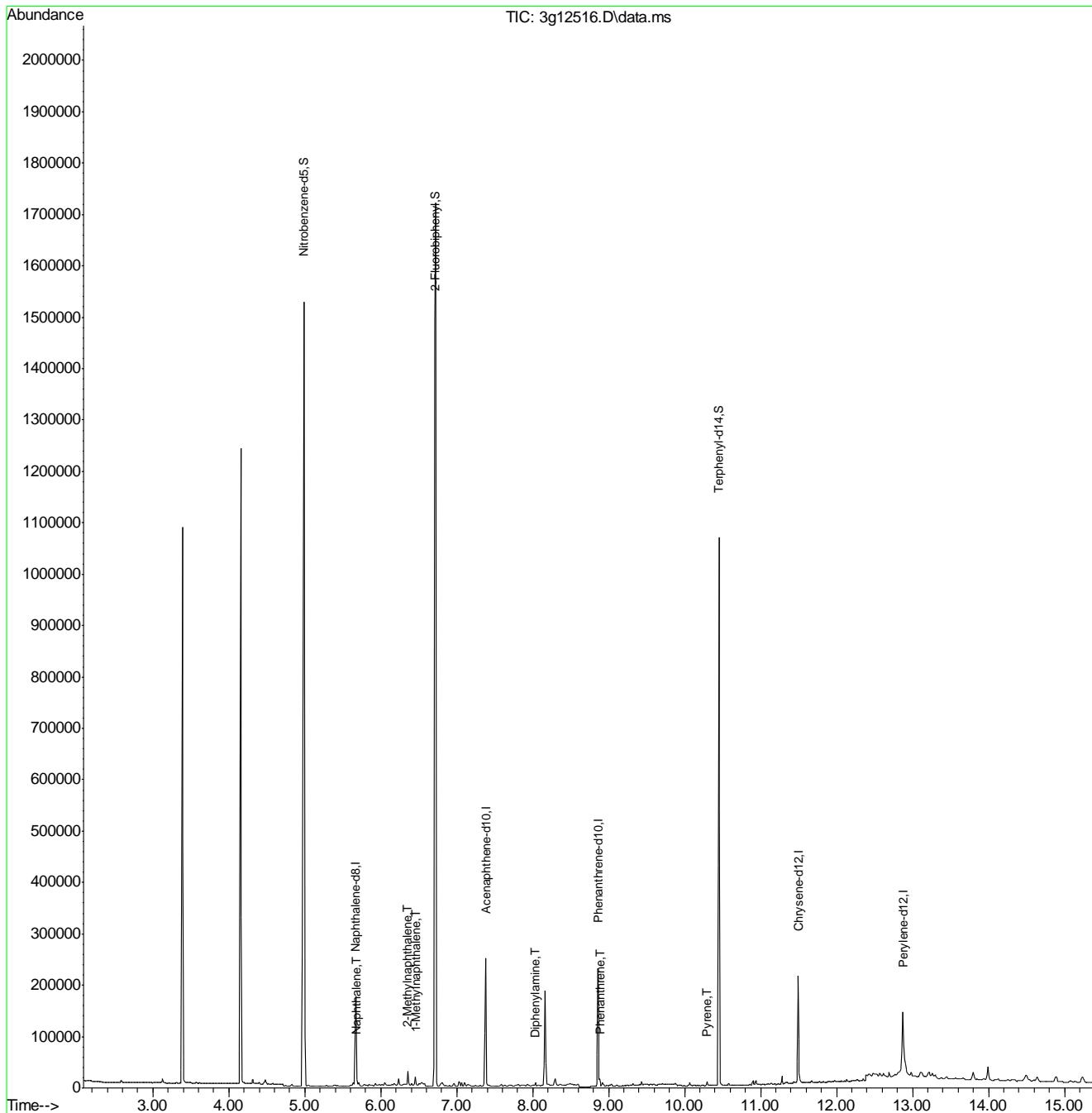
Target Compounds					Qvalue
3) N-Nitrosodimethylamine	2.363	74	25	N.D.	
4) N-Nitrosodi-propylamine	0.000	70	0	N.D. d	
5) Naphthalene	5.683	128	7201	0.1566	ug/mL 93
8) 2-Methylnaphthalene	6.356	142	10319	0.2963	ug/mL 97
9) 1-Methylnaphthalene	6.456	142	6113	0.1771	ug/mL 94
10) Acenaphthylene	7.243	152	473	N.D.	
11) Acenaphthene	7.384	154	461	Below Cal #	20
12) Dibenzofuran	7.585	168	1530	N.D.	
13) Fluorene	7.928	166	1942	N.D.	
14) Diphenylamine	8.034	169	2919	0.0773	ug/mL 74
16) Phenanthrene	8.883	178	8507	0.1266	ug/mL# 75
17) Anthracene	8.938	178	1598	N.D.	
18) Fluoranthene	10.070	202	3235	N.D.	
20) Pyrene	10.292	202	4179	0.0657	ug/mL 93
22) Benzo(a)anthracene	11.483	228	1750	N.D.	
23) Chrysene	11.516	228	2212	N.D.	
25) Benzo(b)fluoranthene	0.000	252	0	N.D. d	
26) Benzo(k)fluoranthene	12.494	252	2527	N.D.	
27) Benzo(a)pyrene	12.820	252	917	N.D.	
28) Indeno(1,2,3-cd)pyrene	14.082	276	618	N.D.	
29) Dibenz(a,h)anthracene	14.103	278	483	N.D.	
30) Benzo(g,h,i)perylene	14.440	276	1738	N.D.	

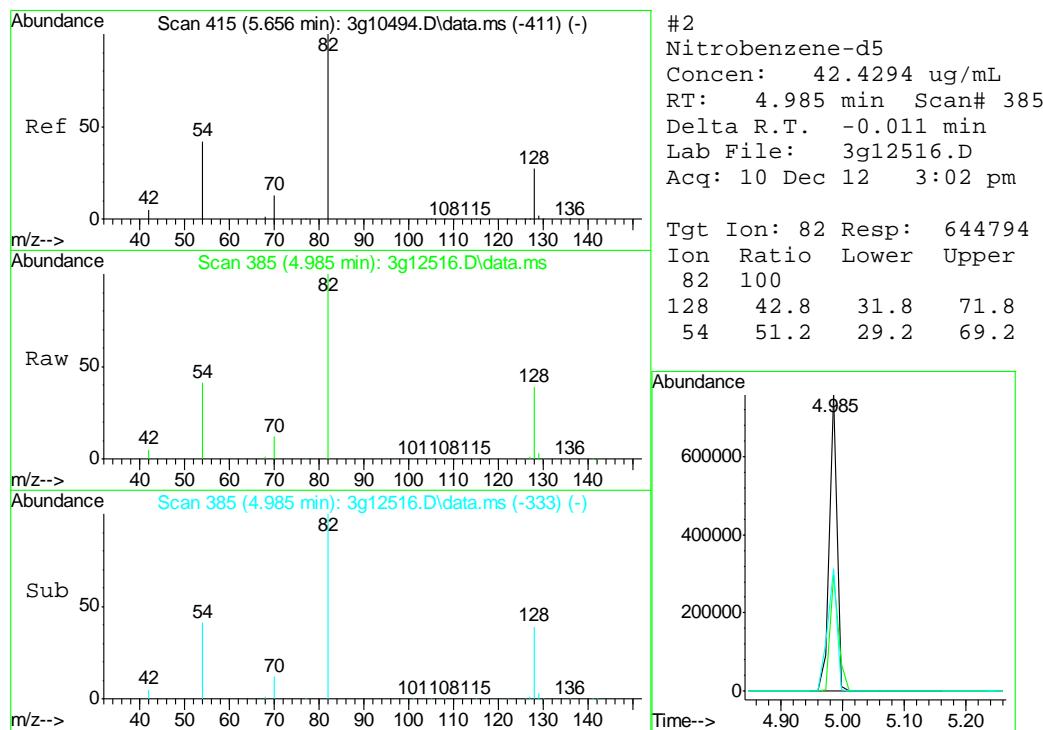
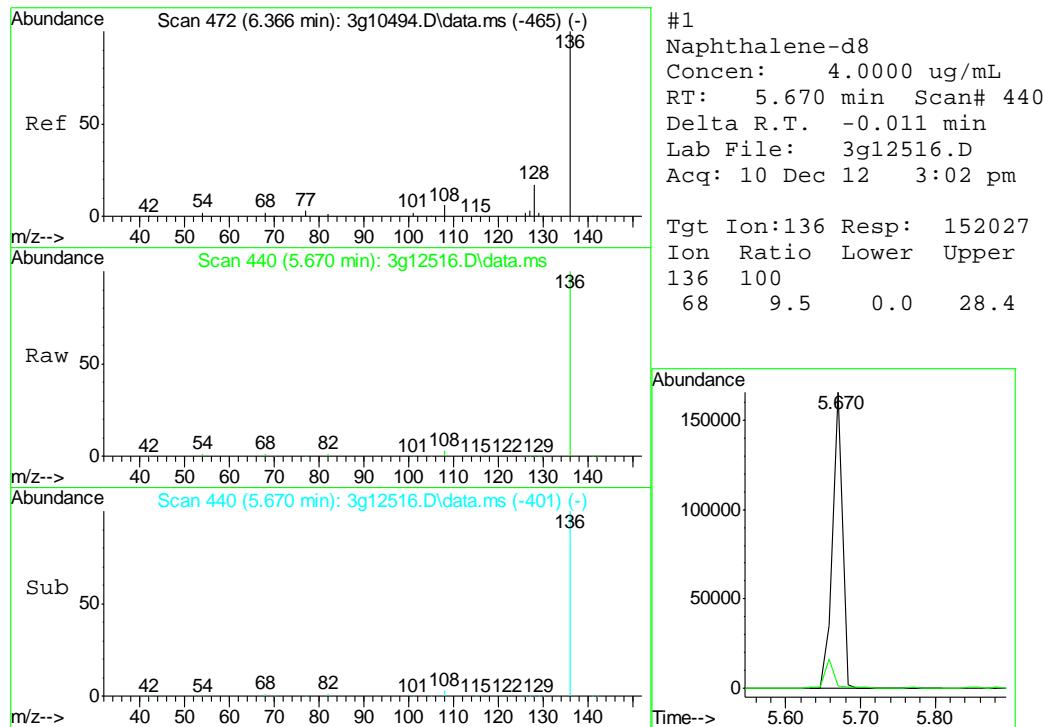
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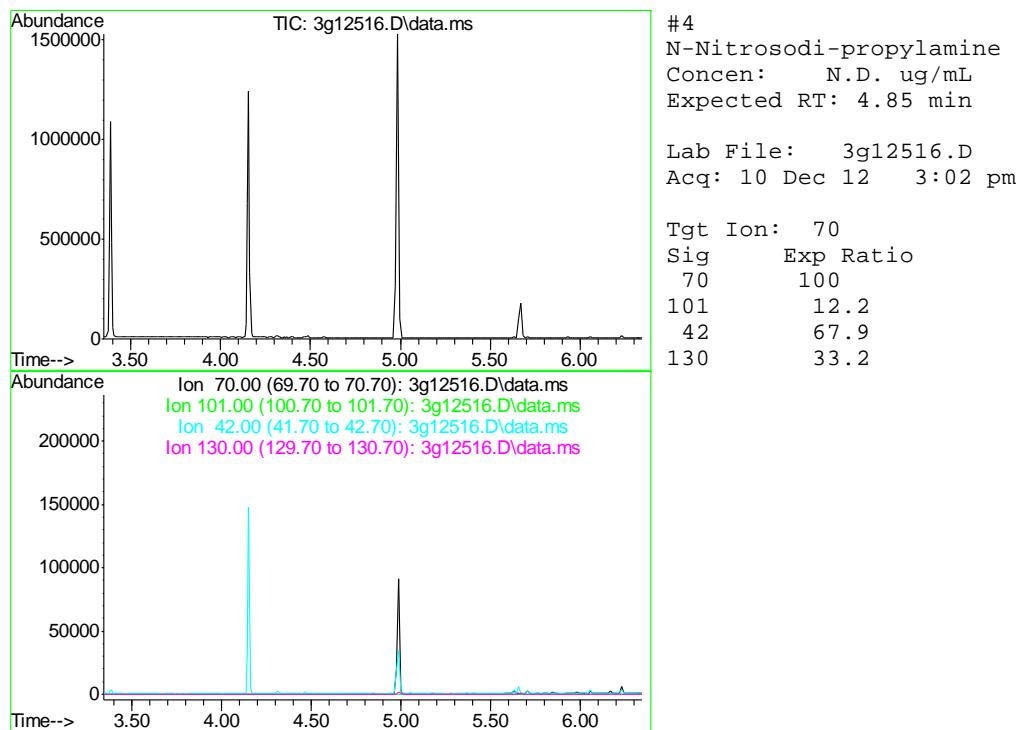
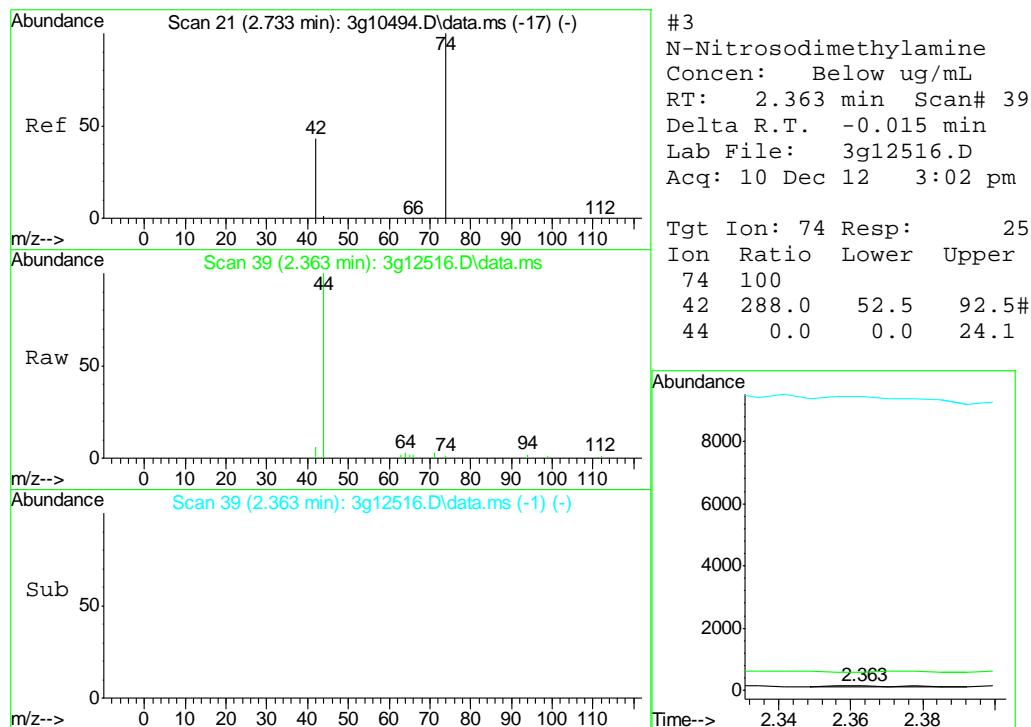
## Quantitation Report (QT Reviewed)

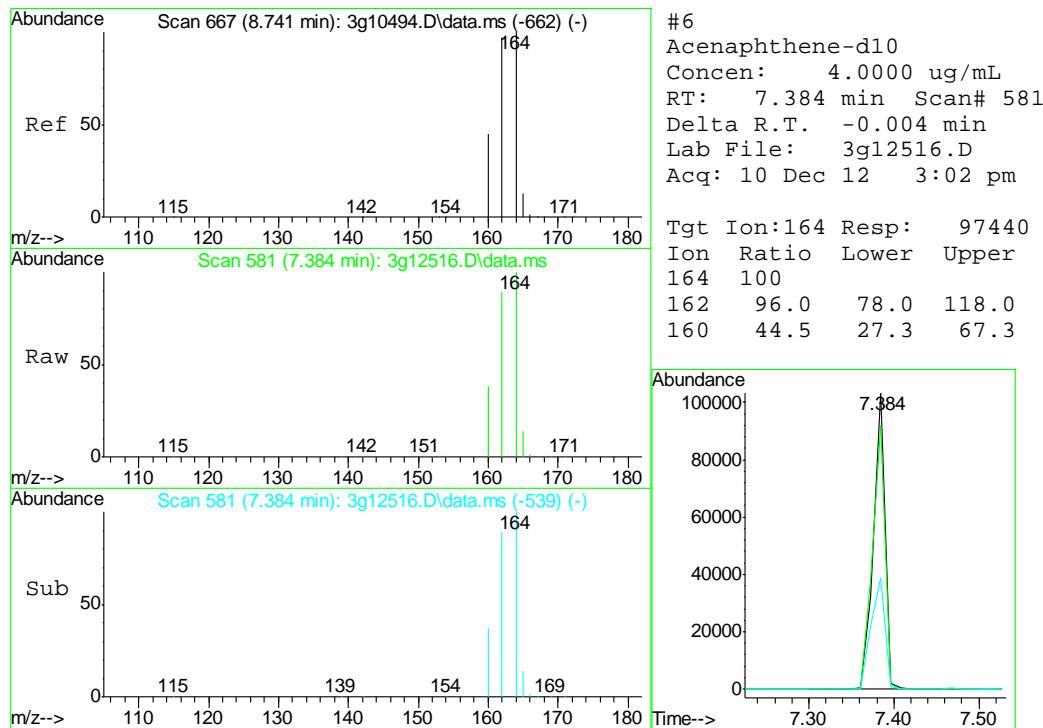
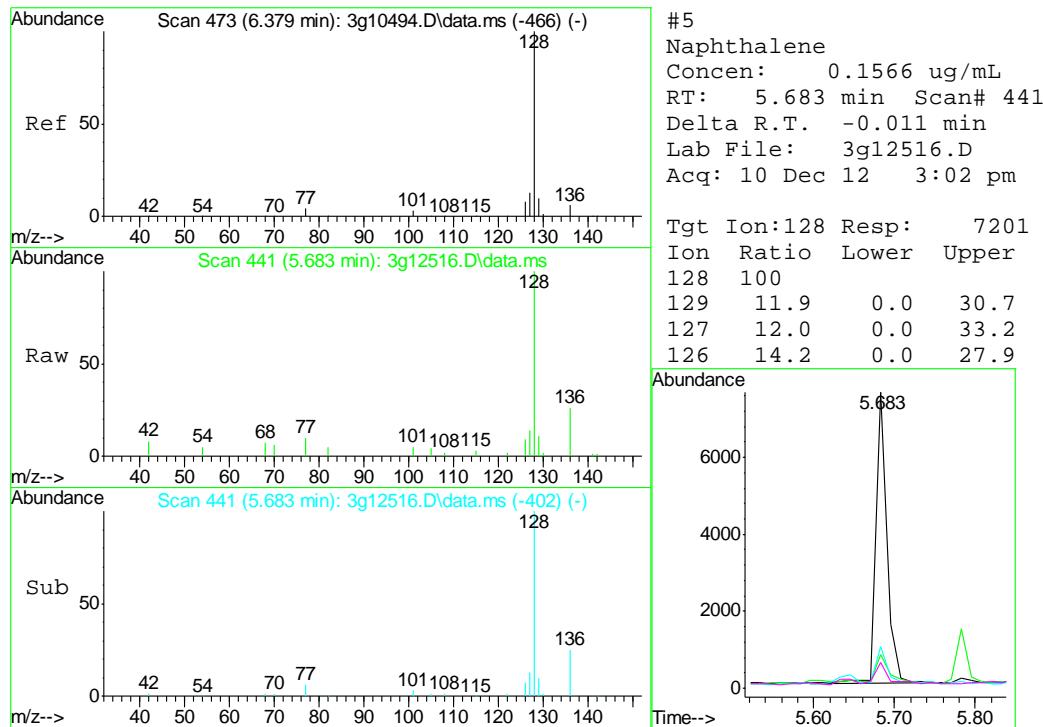
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 Sample : D41506-1  
 Misc : OP7075,E3G593,30.16,,,1,1  
 ALS Vial : 12 Sample Multiplier: 1

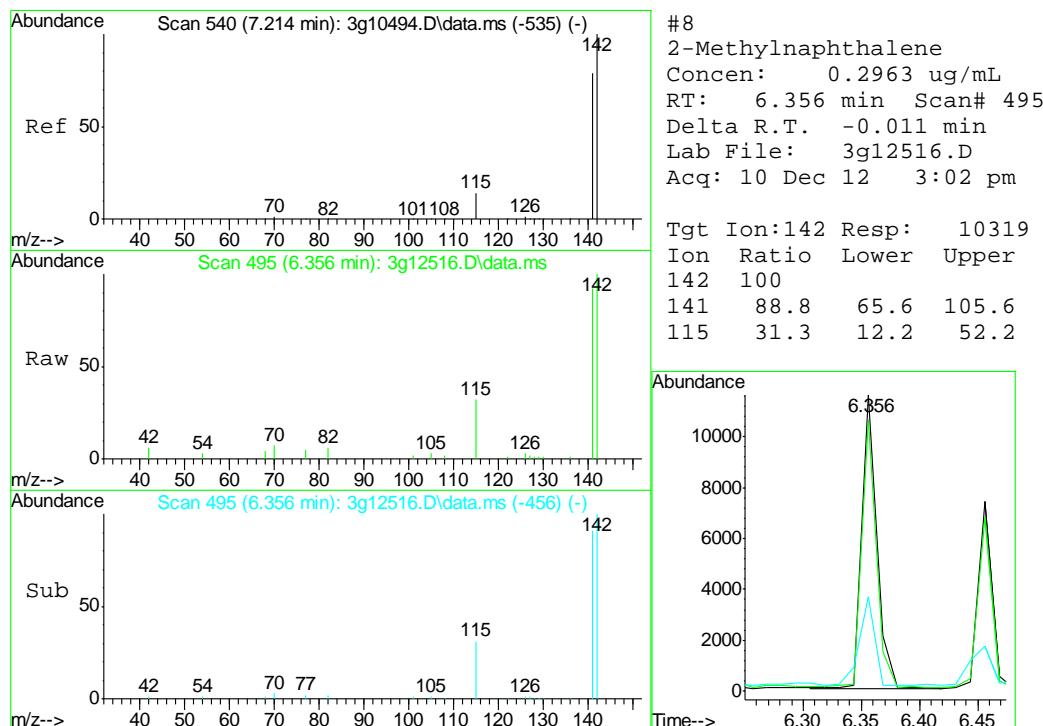
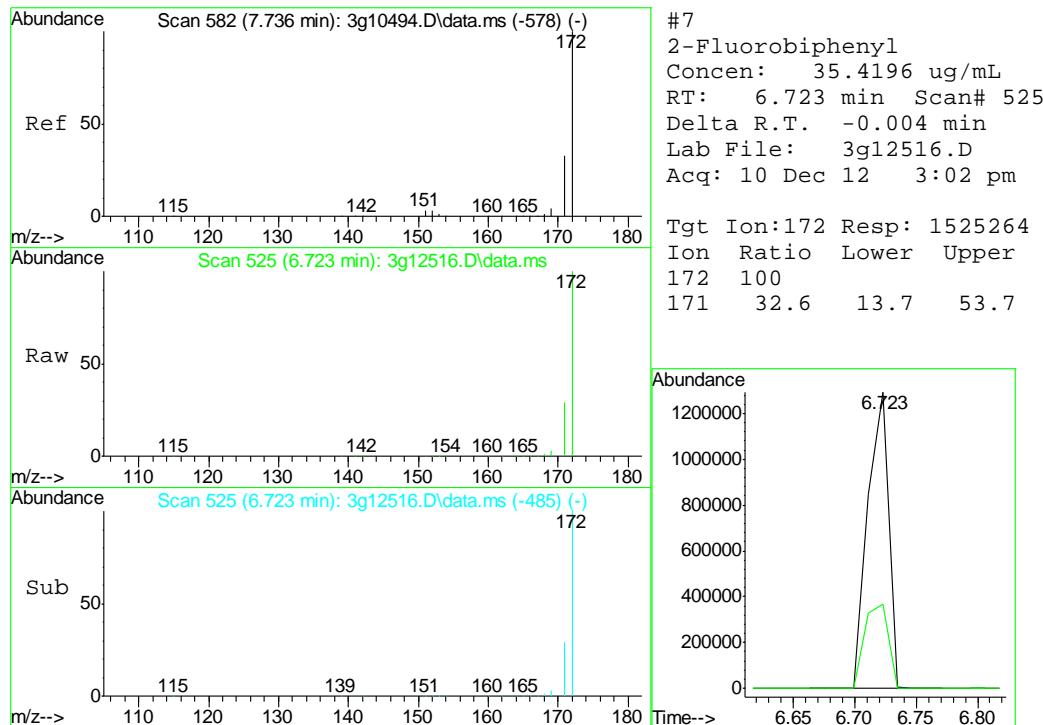
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 Quant Title : PAHSIM BASE  
 QLast Update : Tue Dec 04 08:50:28 2012  
 Response via : Initial Calibration

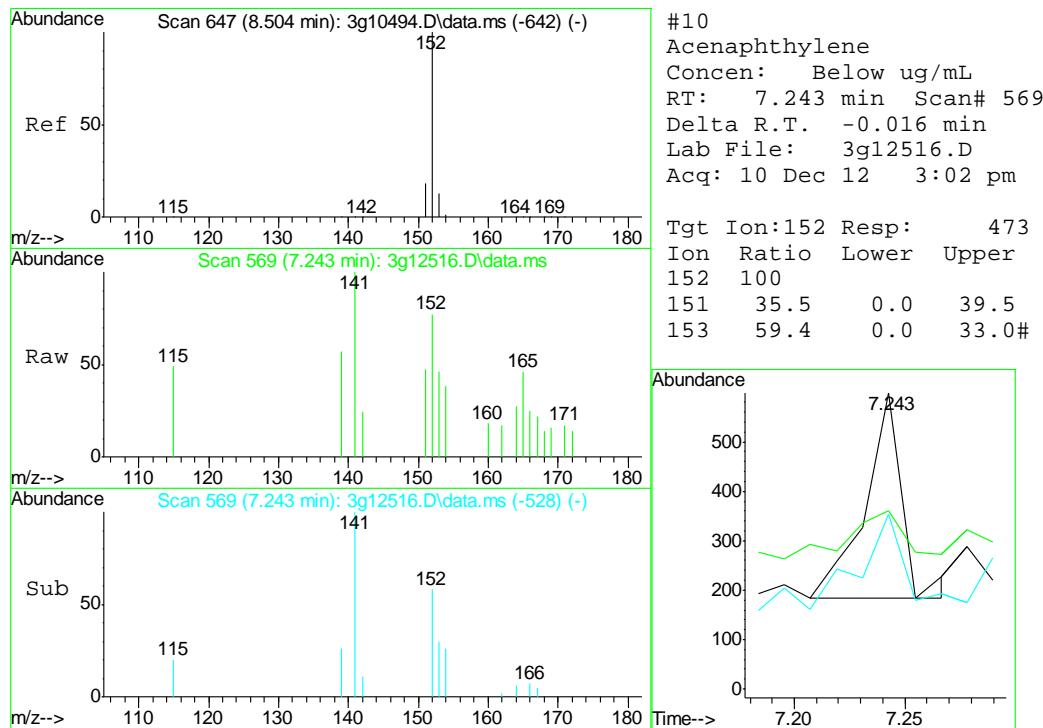
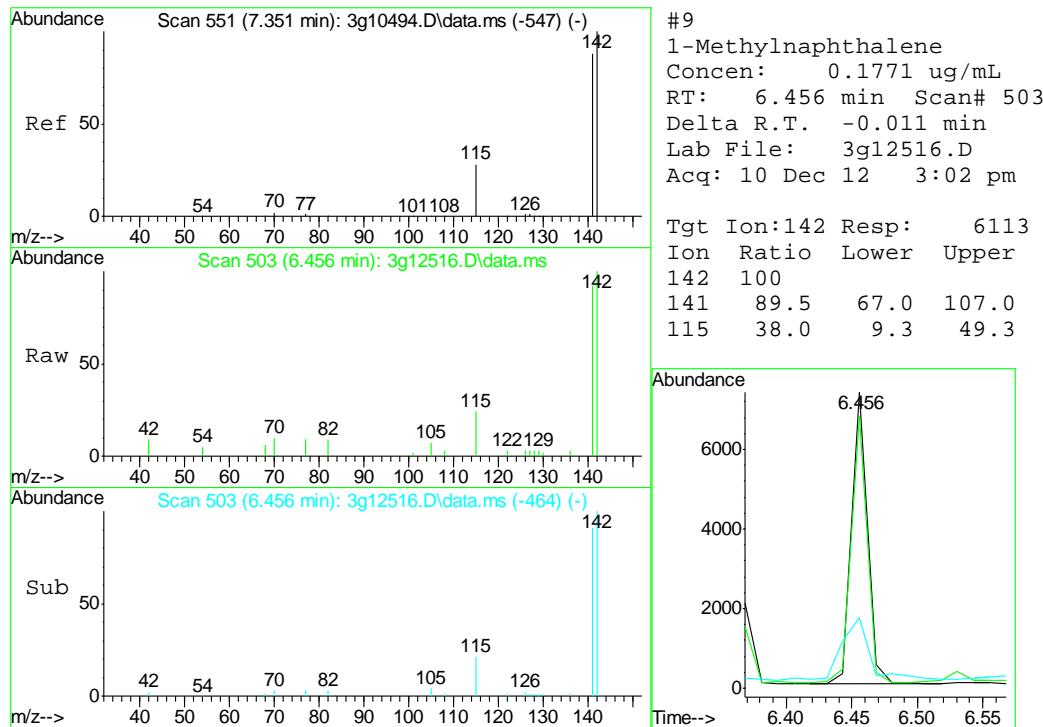


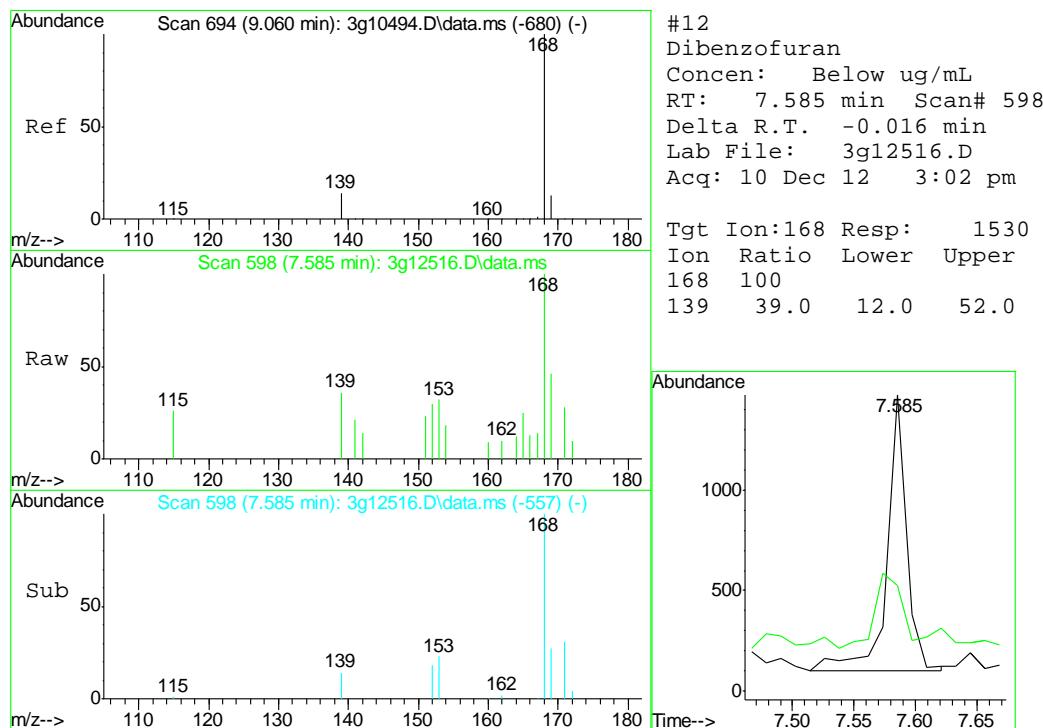
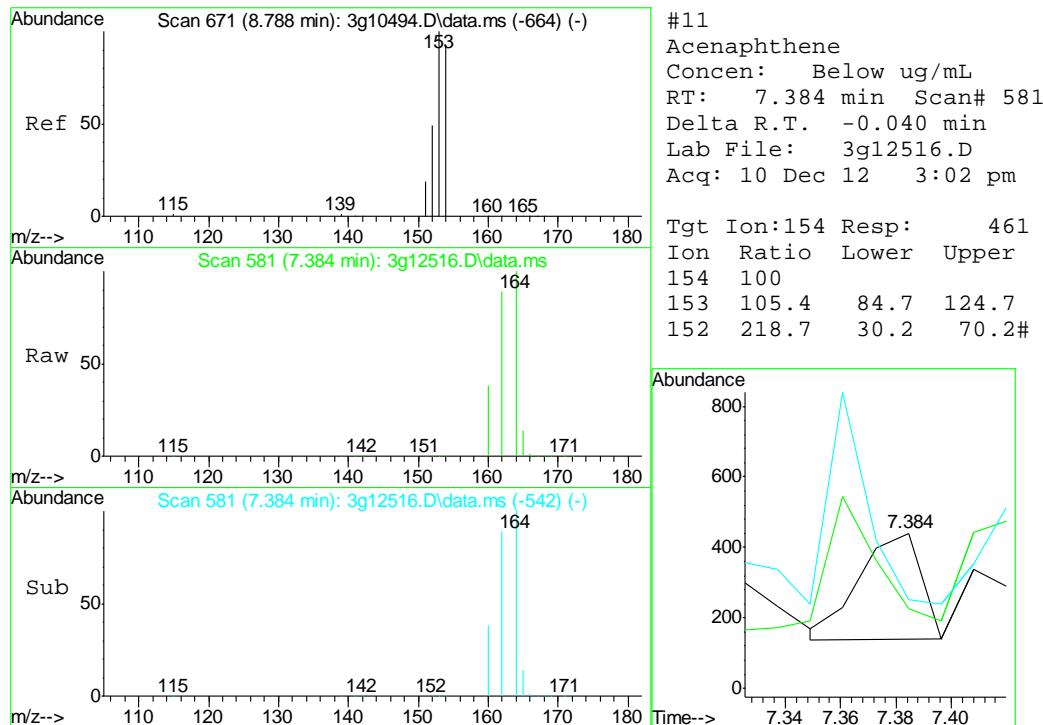


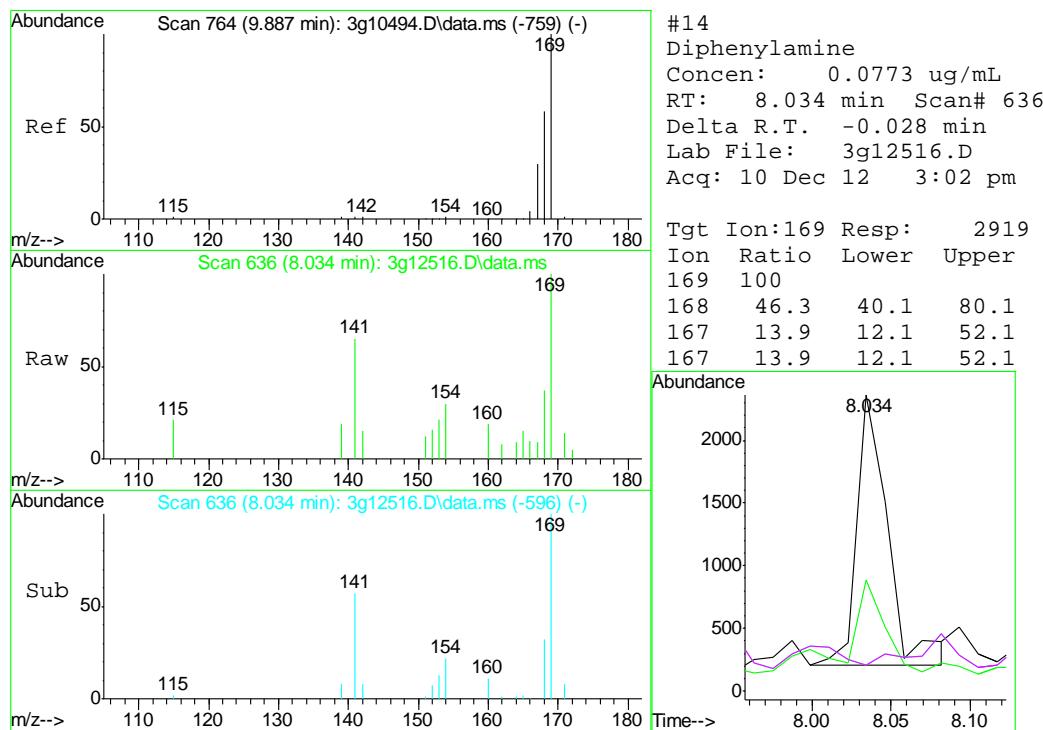
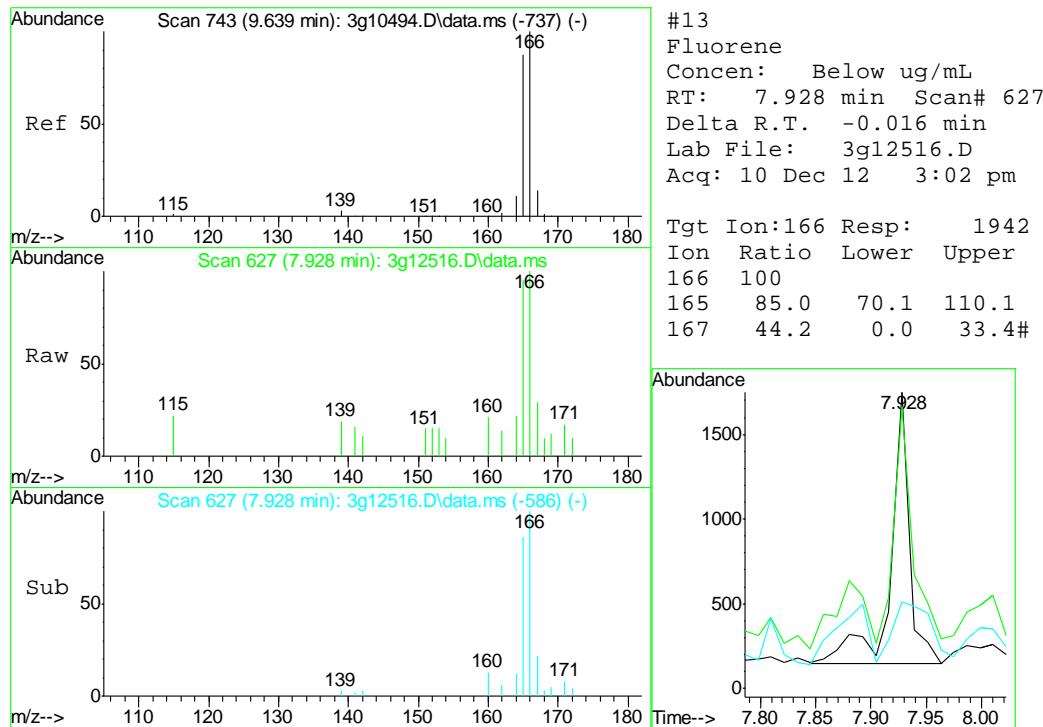


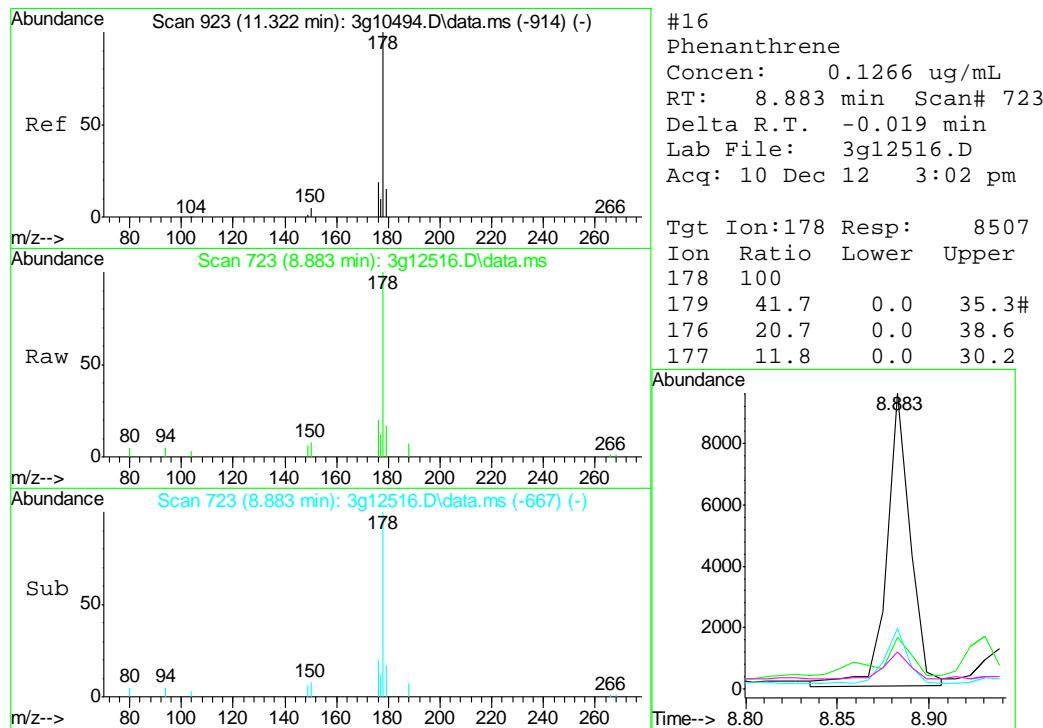
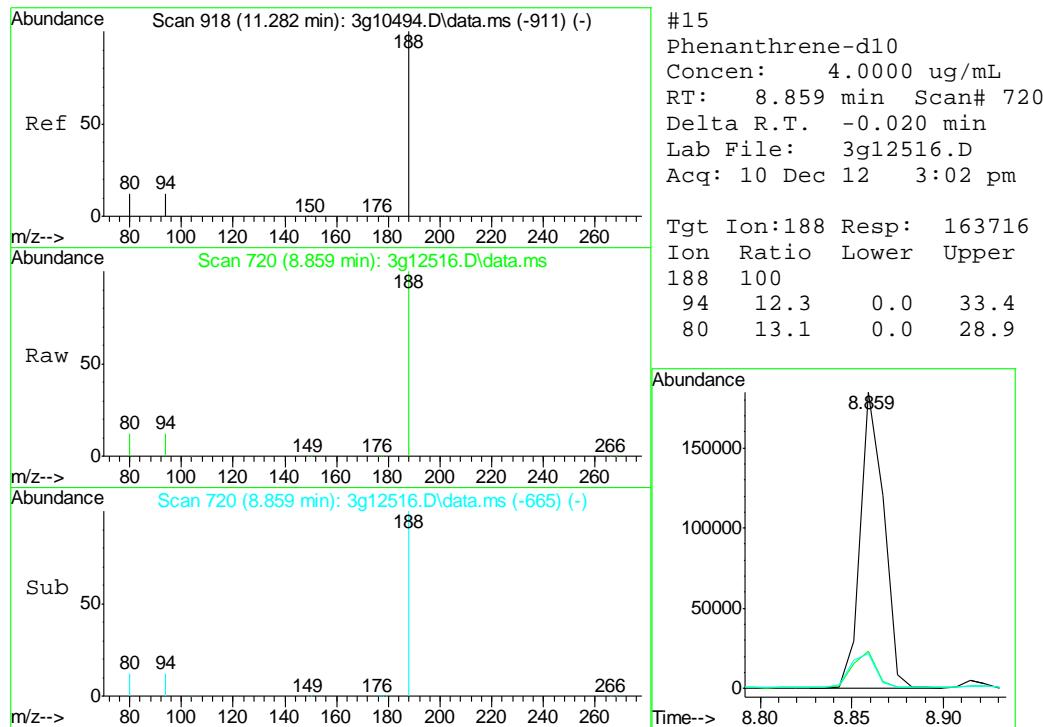


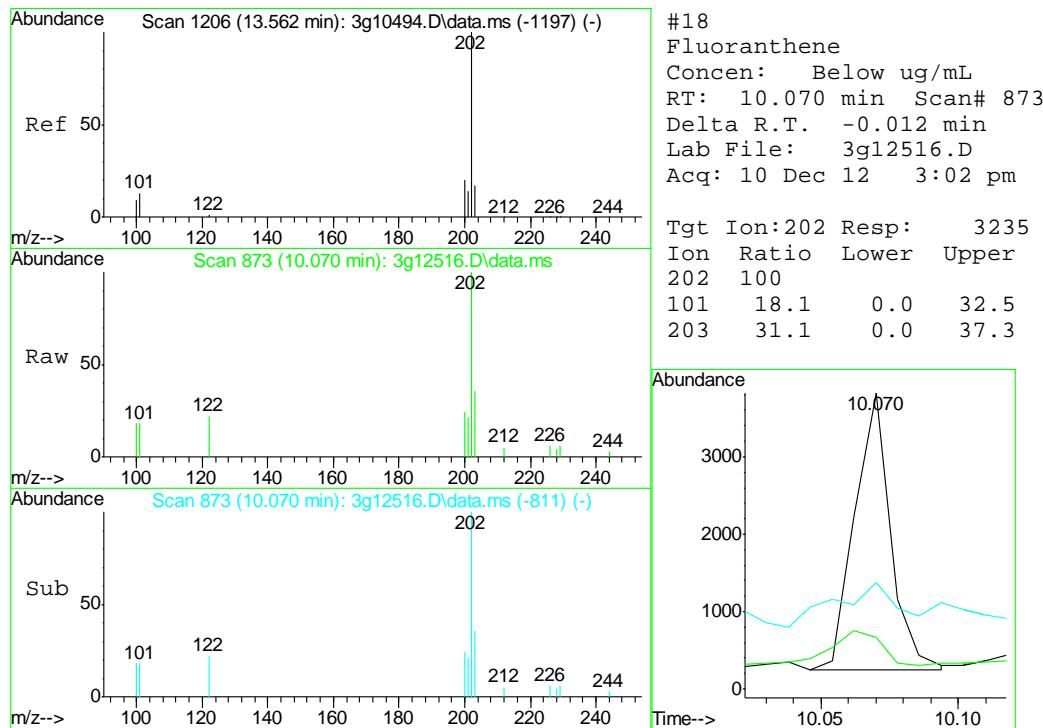
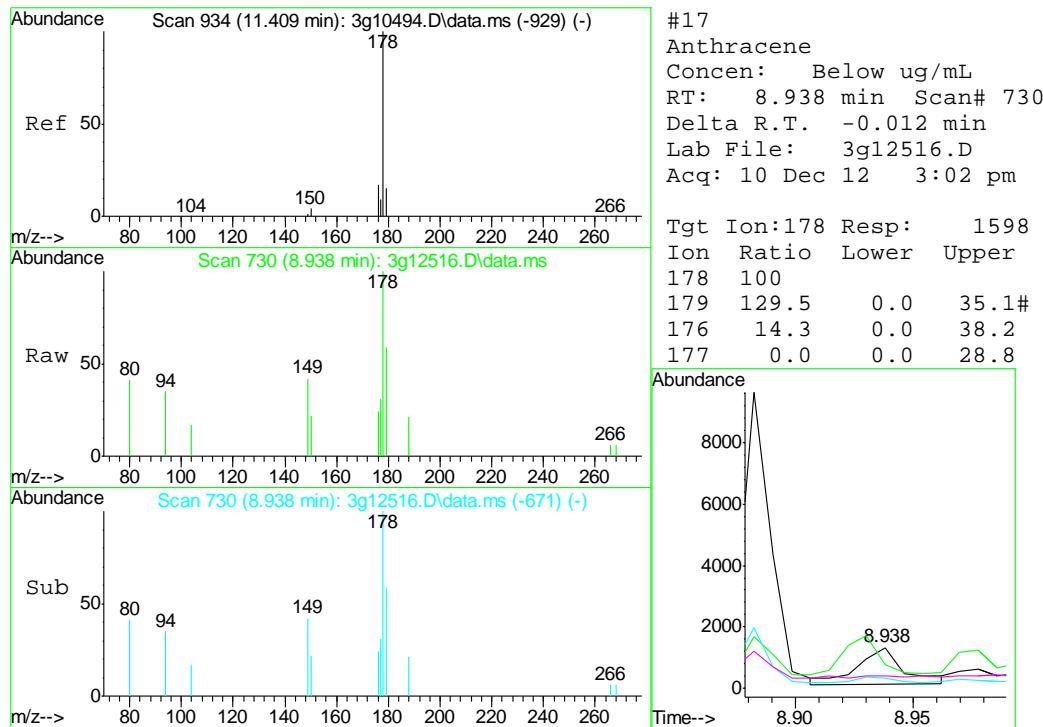


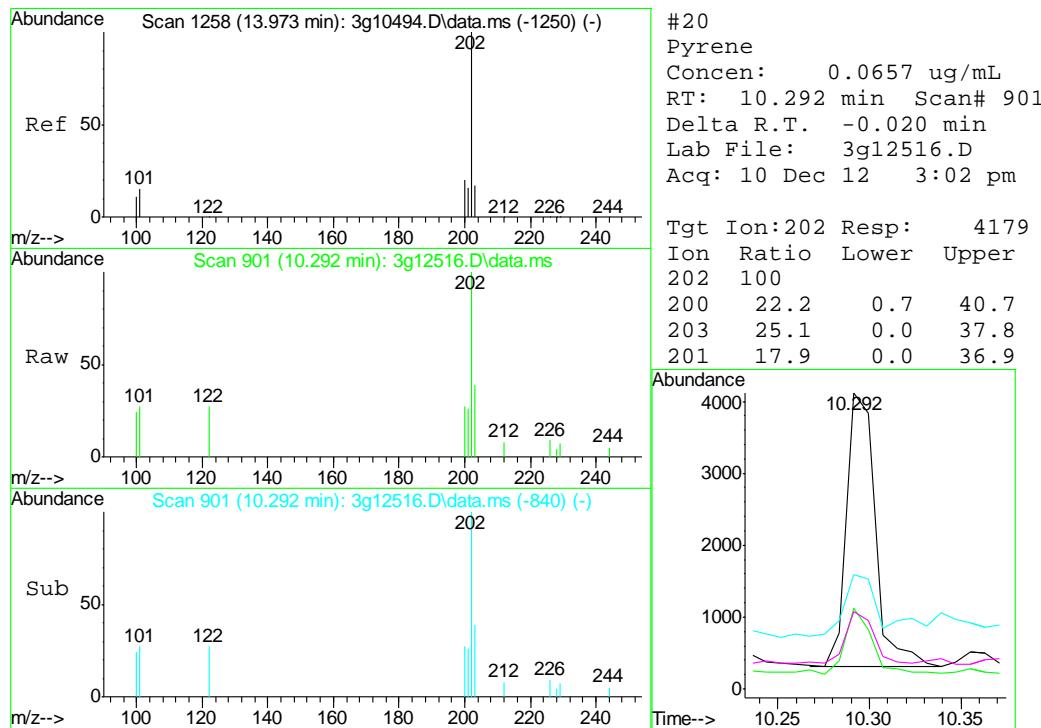
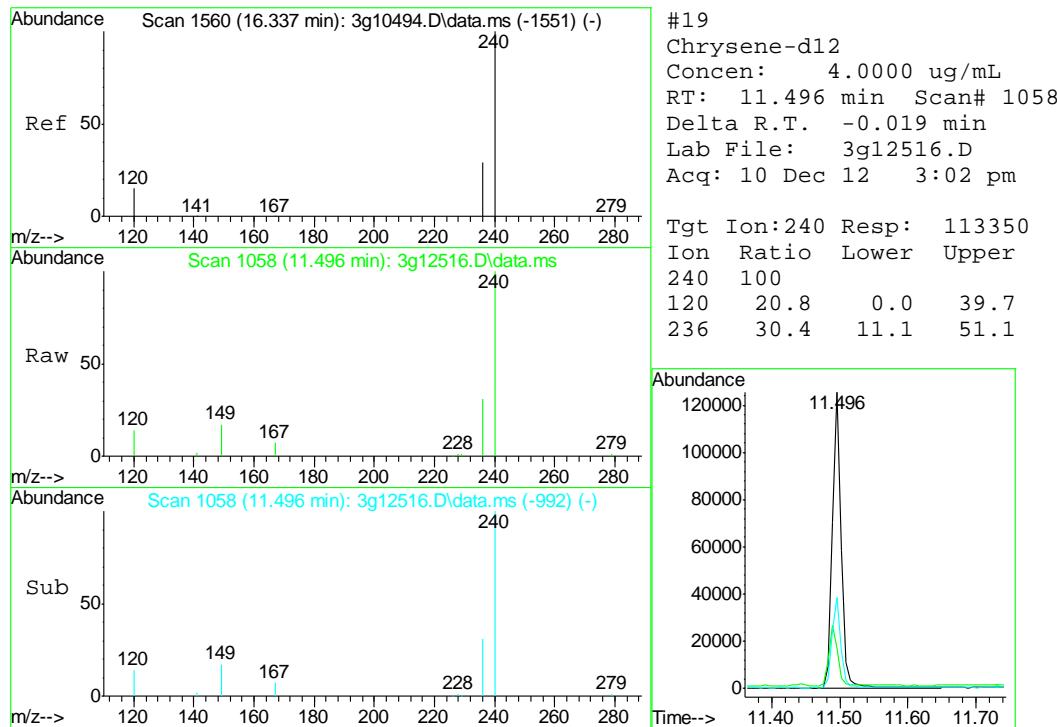


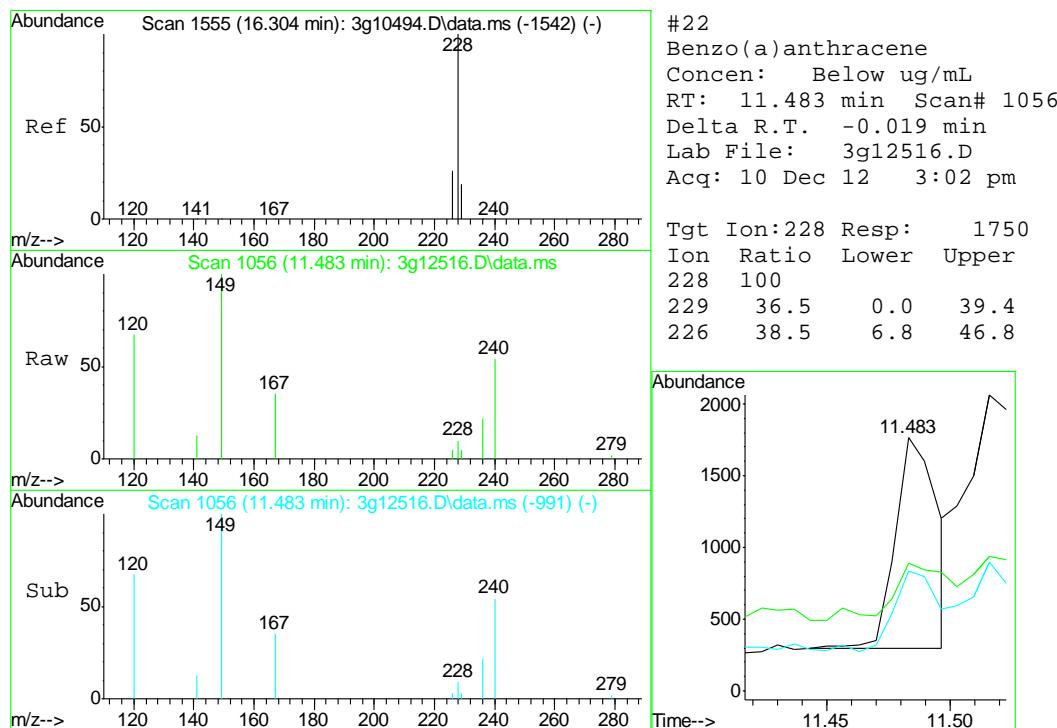
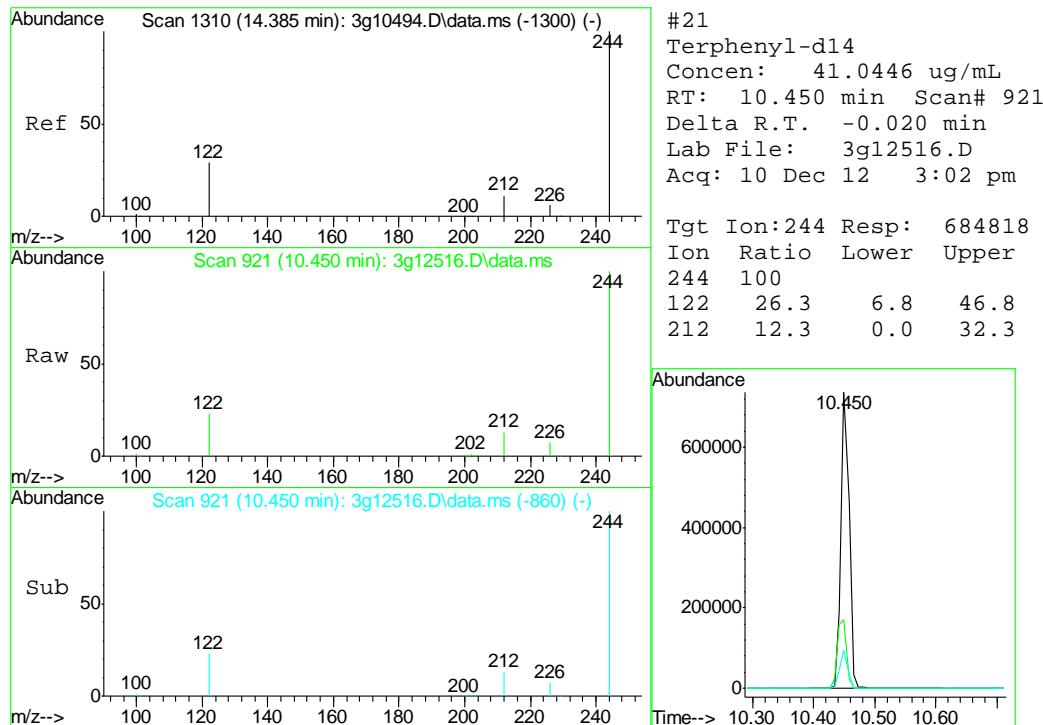


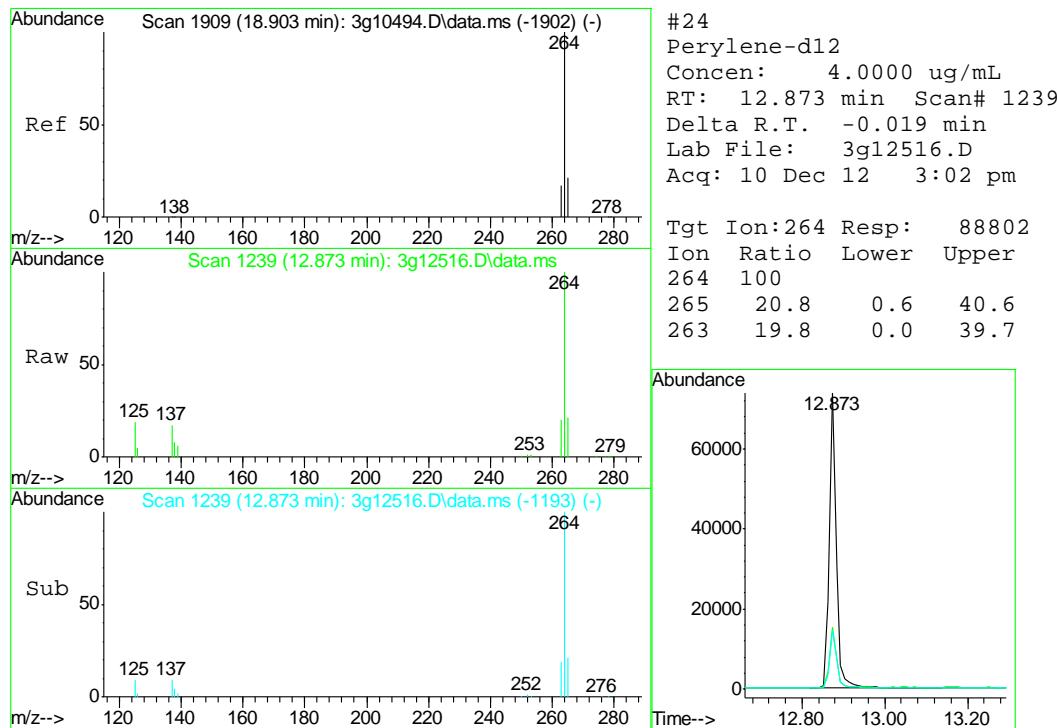
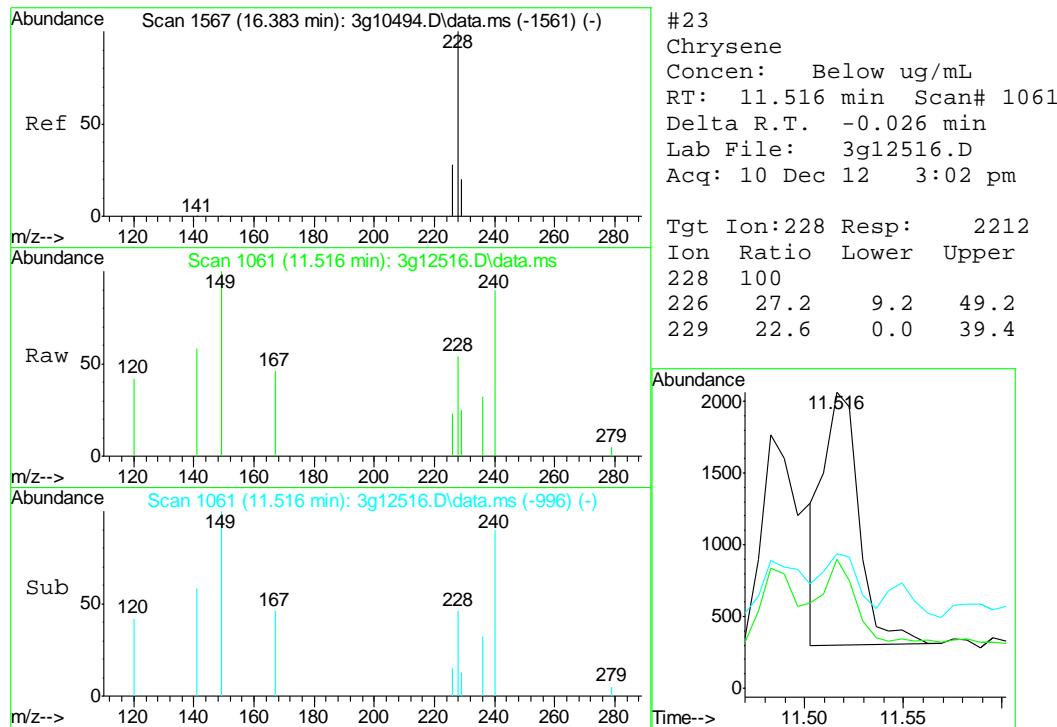


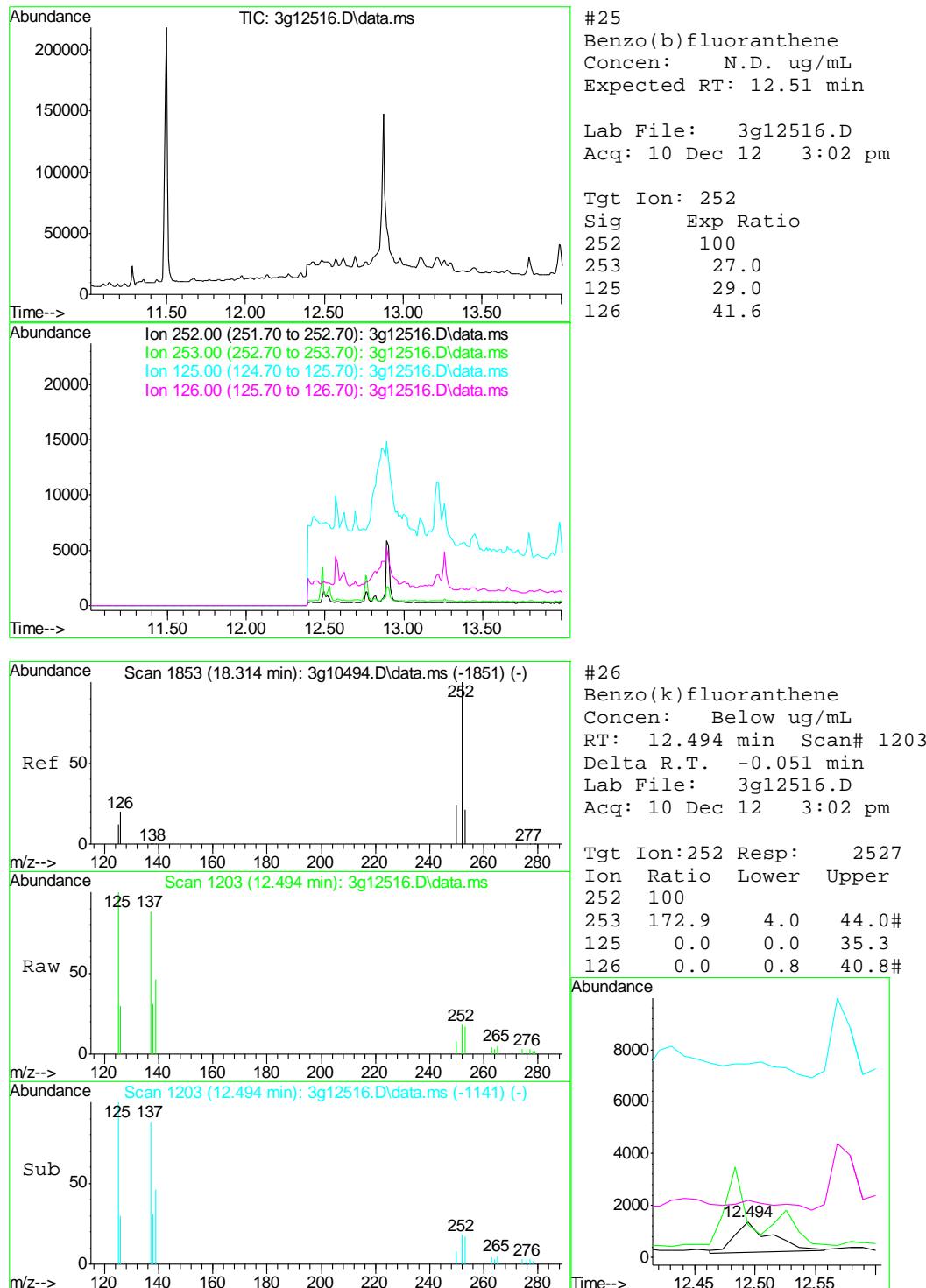


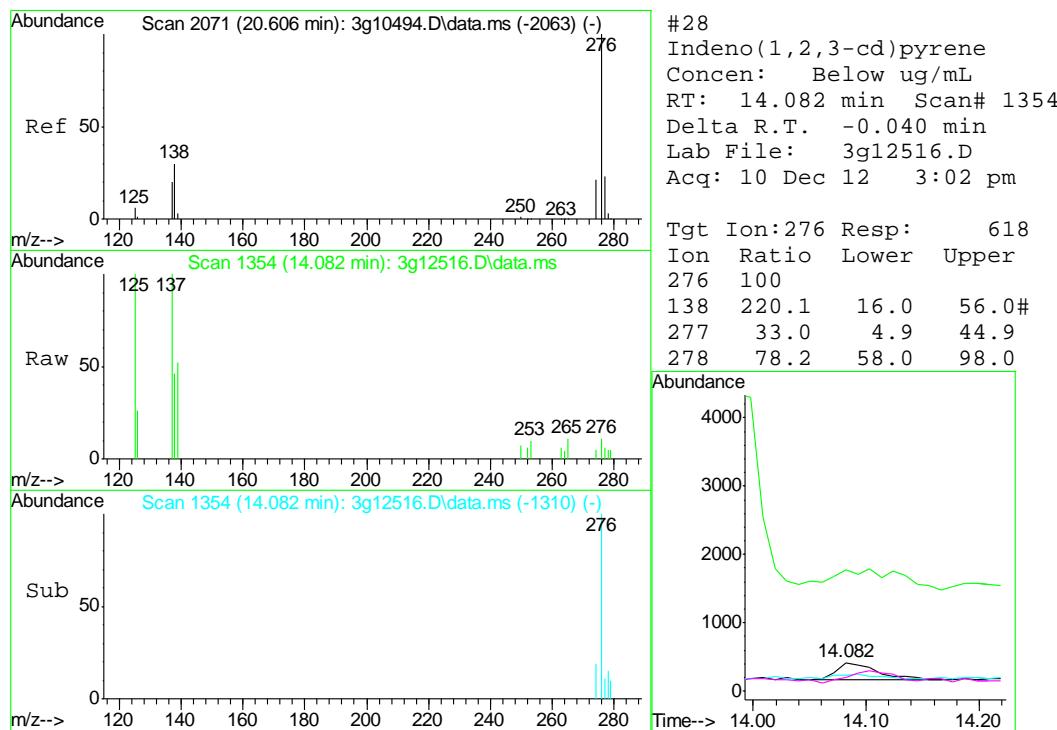
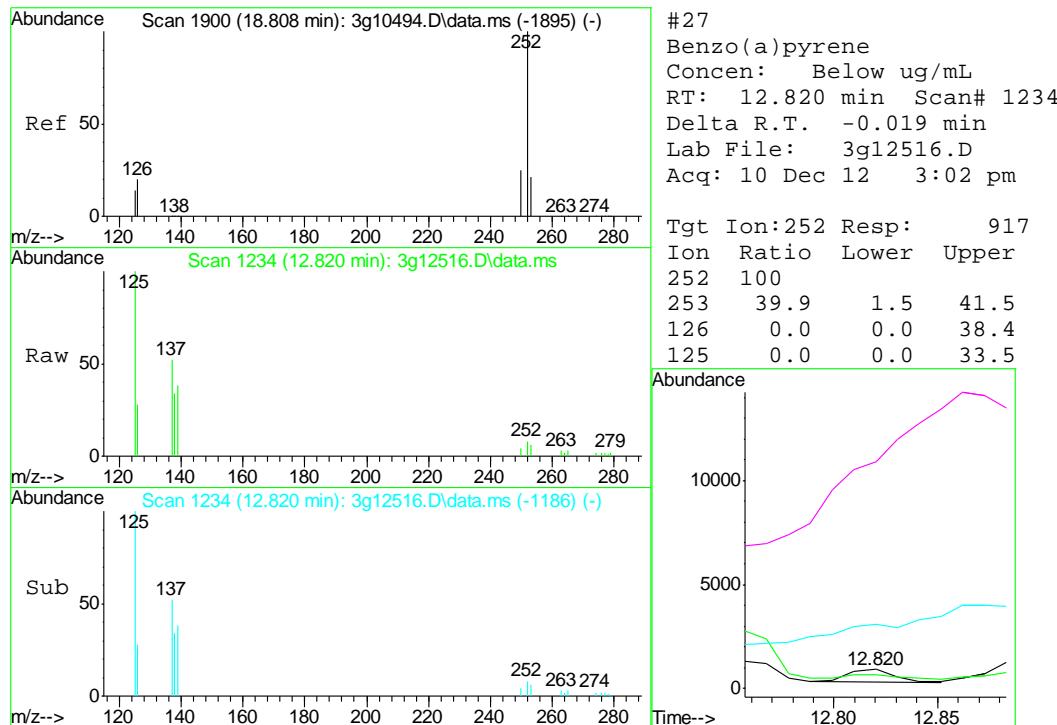


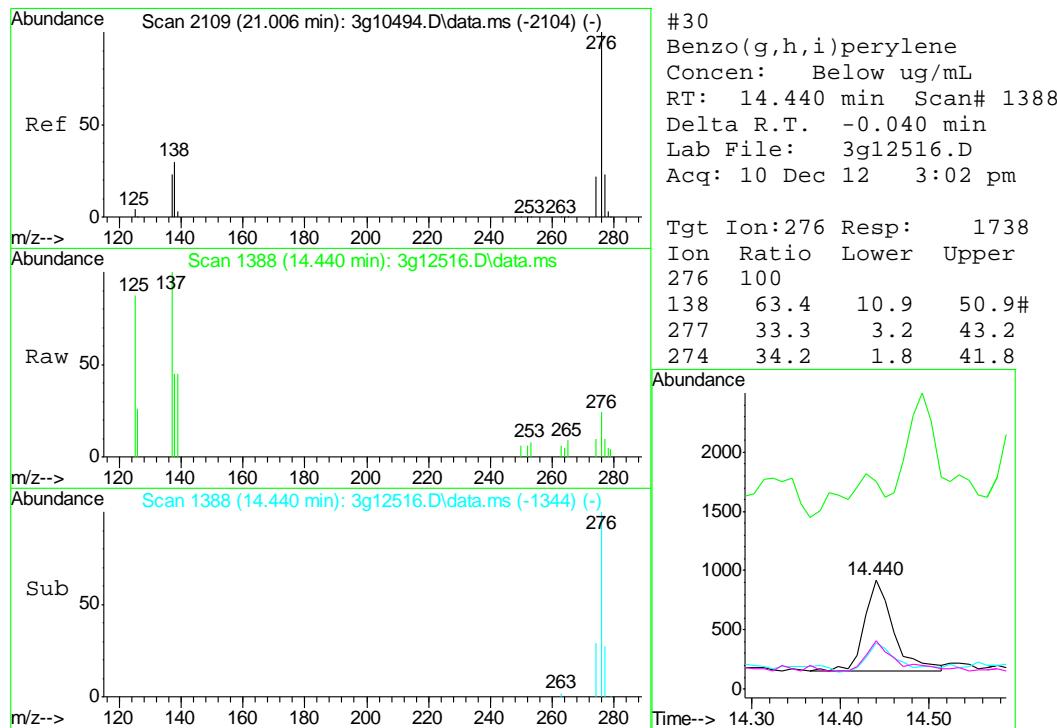
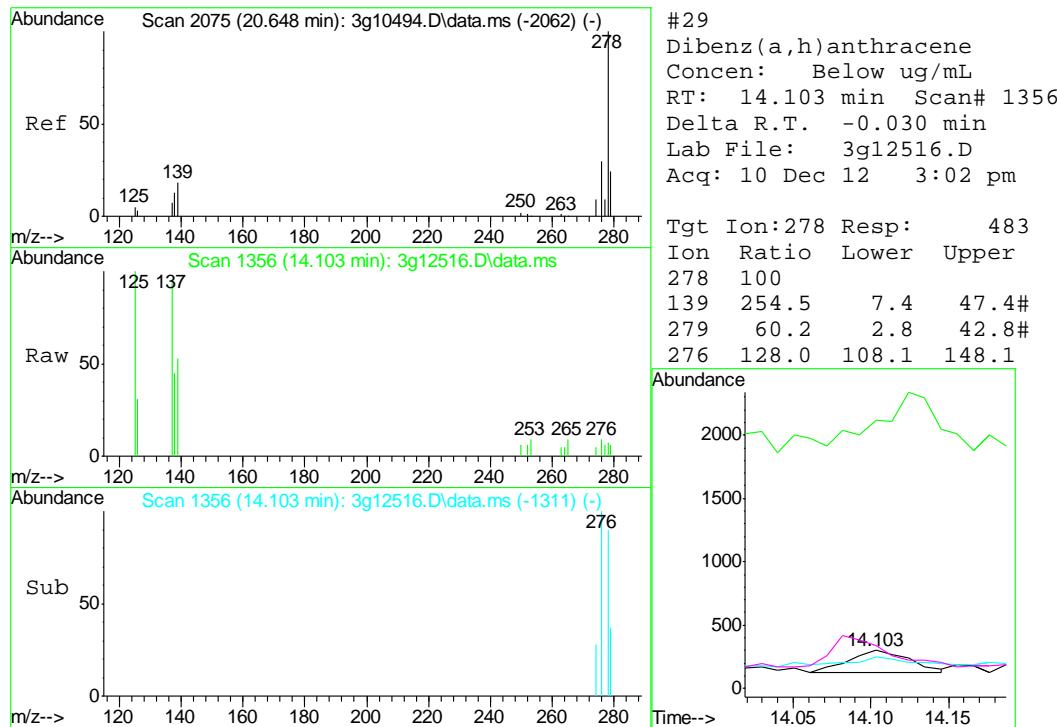












## Quantitation Report (QT Reviewed)

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 Data File : 3g12508.D  
 Acq On : 10 Dec 2012 11:52 am  
 Operator : DONC  
 Sample : OP7075-MB  
 Misc : OP7075,E3G593,30.00,,,1,1  
 ALS Vial : 4 Sample Multiplier: 1

Quant Time: Dec 10 13:58:13 2012  
 Quant Method : C:\msdchem\1\METHODS\SIMPE3G586.M  
 Quant Title : PAHSIM BASE  
 QLast Update : Tue Dec 04 08:50:28 2012  
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Naphthalene-d8	5.670	136	153335	4.0000	ug/mL	-0.01
6) Acenaphthene-d10	7.385	164	92403	4.0000	ug/mL	0.00
15) Phenanthrene-d10	8.867	188	160356	4.0000	ug/mL	-0.01
19) Chrysene-d12	11.503	240	115791	4.0000	ug/mL	-0.01
24) Perylene-d12	12.883	264	94737	4.0000	ug/mL	0.00

System Monitoring Compounds						
2) Nitrobenzene-d5	4.985	82	677706	44.2147	ug/mL	-0.01
Spiked Amount	50.000	Range	25 - 135	Recovery	=	88.42%
7) 2-Fluorobiphenyl	6.723	172	1592975	39.4888	ug/mL	0.00
Spiked Amount	50.000	Range	25 - 135	Recovery	=	78.98%
21) Terphenyl-d14	10.458	244	774374	45.4337	ug/mL	-0.01
Spiked Amount	50.000	Range	25 - 135	Recovery	=	90.86%

Target Compounds					Qvalue
3) N-Nitrosodimethylamine	2.334	74	50	N.D.	
4) N-Nitrosodi-propylamine	0.000	70	0	N.D. d	
5) Naphthalene	5.683	128	491	N.D.	
8) 2-Methylnaphthalene	6.356	142	179	N.D.	
9) 1-Methylnaphthalene	6.456	142	126	N.D.	
10) Acenaphthylene	7.243	152	198	N.D.	
11) Acenaphthene	7.113	154	75	Below Cal	87
12) Dibenzofuran	7.585	168	152	N.D.	
13) Fluorene	0.000	166	0	N.D. d	
14) Diphenylamine	0.000	169	0	N.D. d	
16) Phenanthrene	8.891	178	540	N.D.	
17) Anthracene	8.938	178	357	N.D.	
18) Fluoranthene	10.070	202	778	N.D.	
20) Pyrene	10.299	202	864	N.D.	
22) Benzo(a)anthracene	11.496	228	1148	N.D.	
23) Chrysene	11.523	228	724	N.D.	
25) Benzo(b)fluoranthene	12.494	252	1411	N.D.	
26) Benzo(k)fluoranthene	12.494	252	1411	N.D.	
27) Benzo(a)pyrene	12.820	252	509	N.D.	
28) Indeno(1,2,3-cd)pyrene	14.093	276	459	N.D.	
29) Dibenz(a,h)anthracene	14.114	278	376	N.D.	
30) Benzo(g,h,i)perylene	14.450	276	479	N.D.	

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

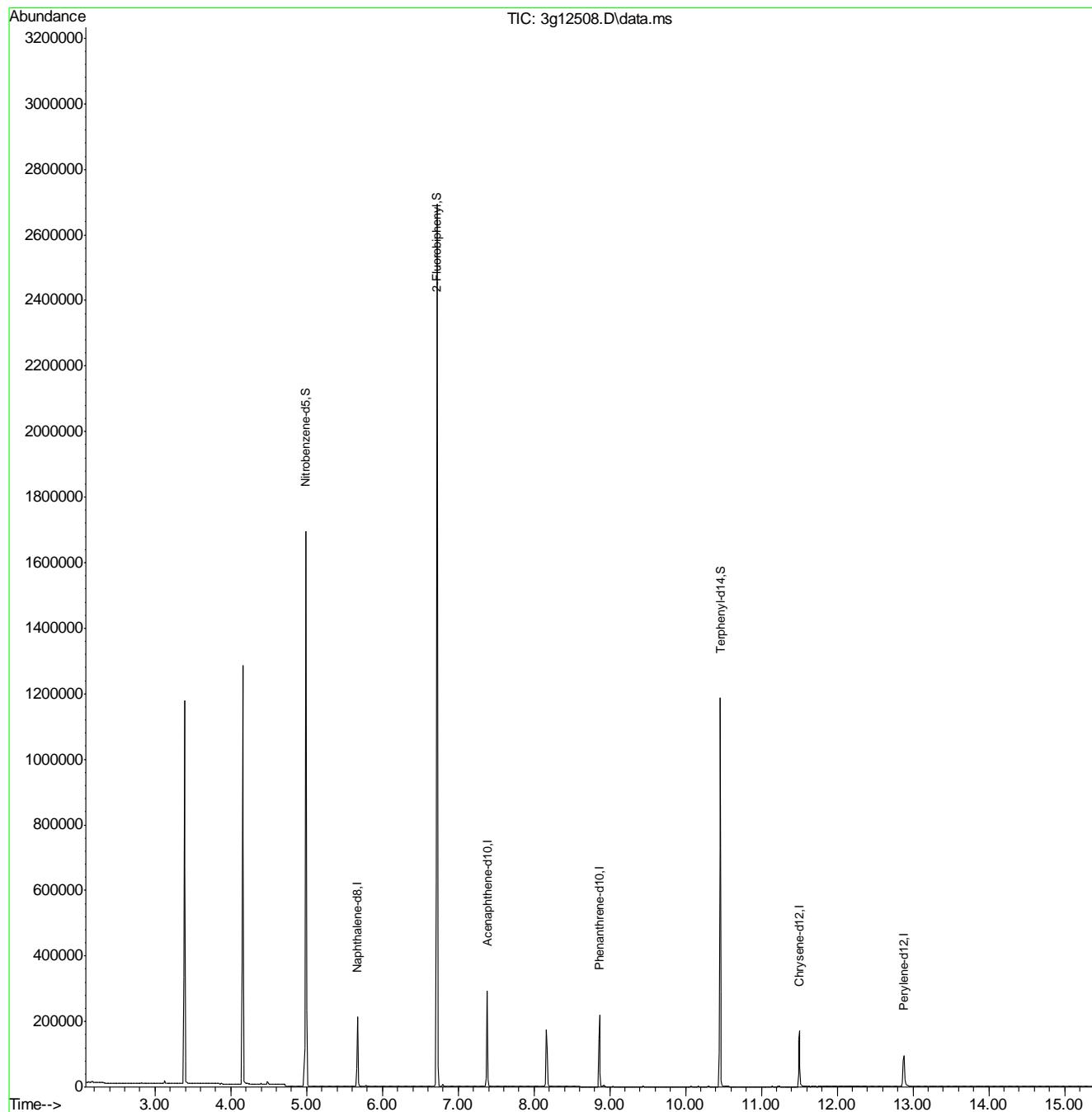
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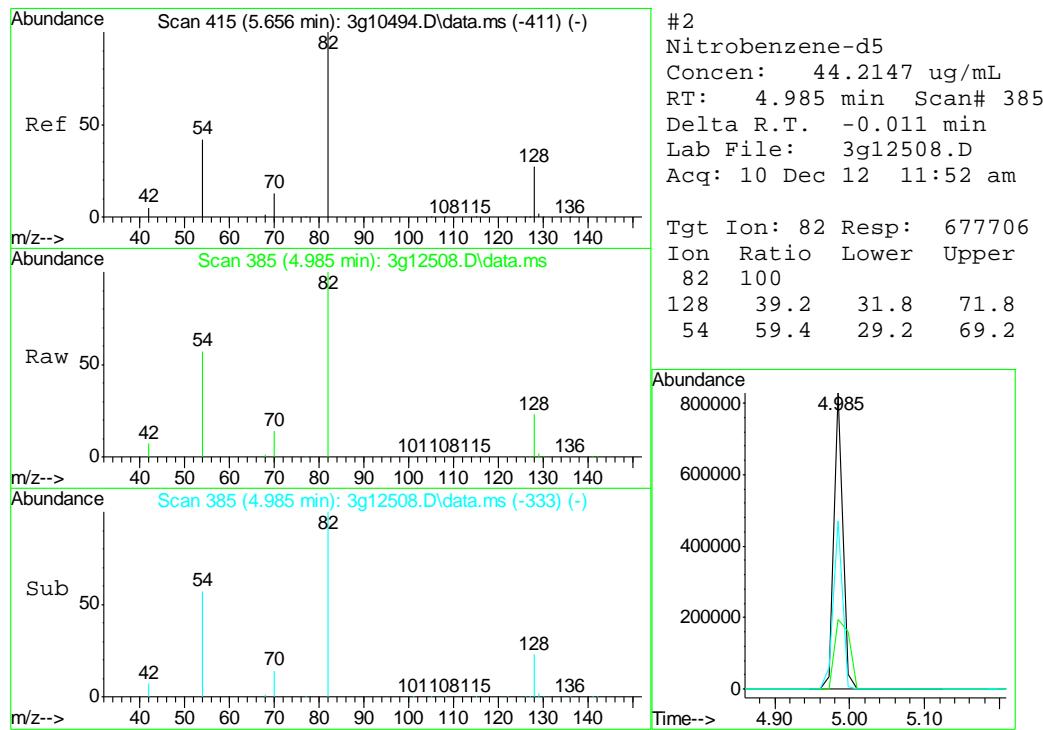
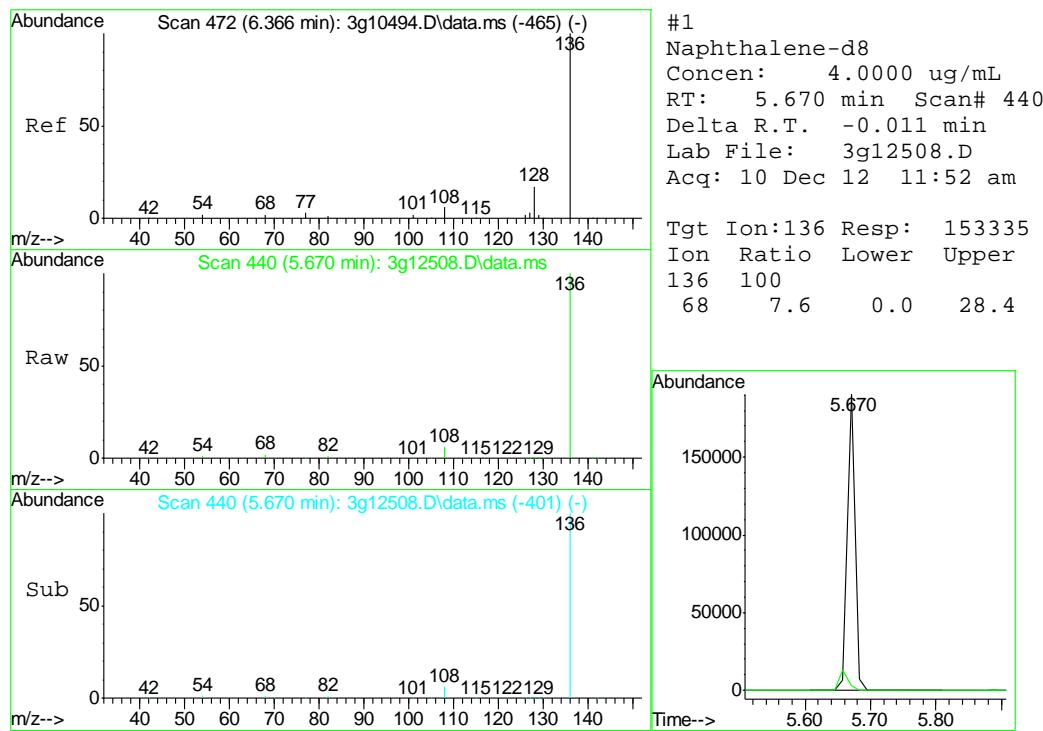
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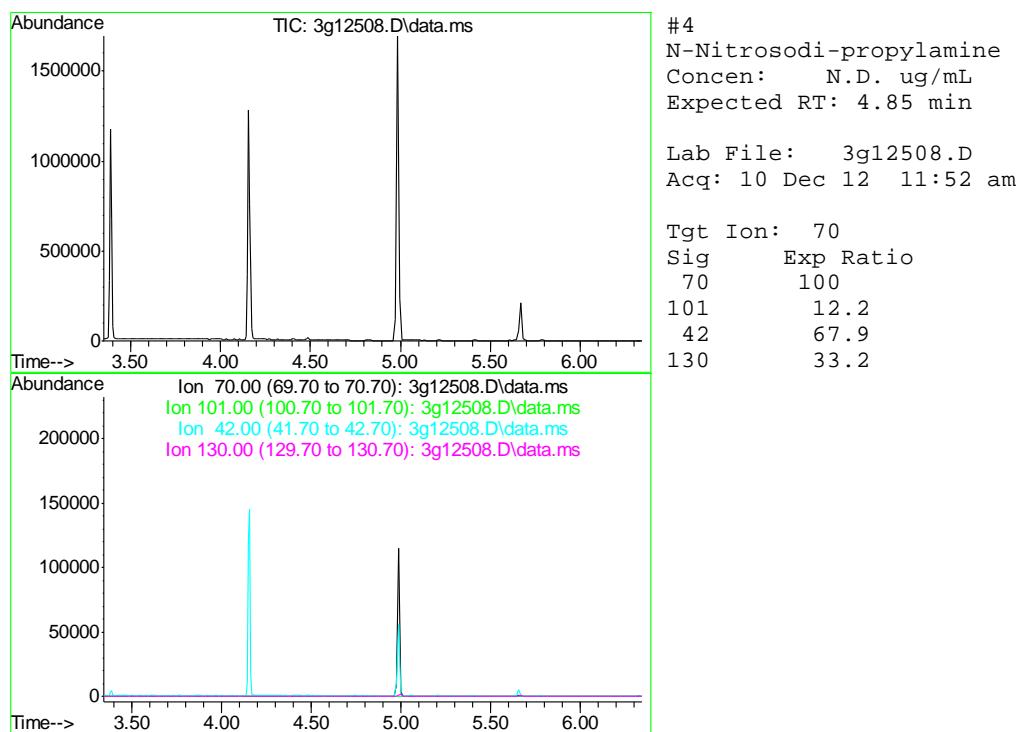
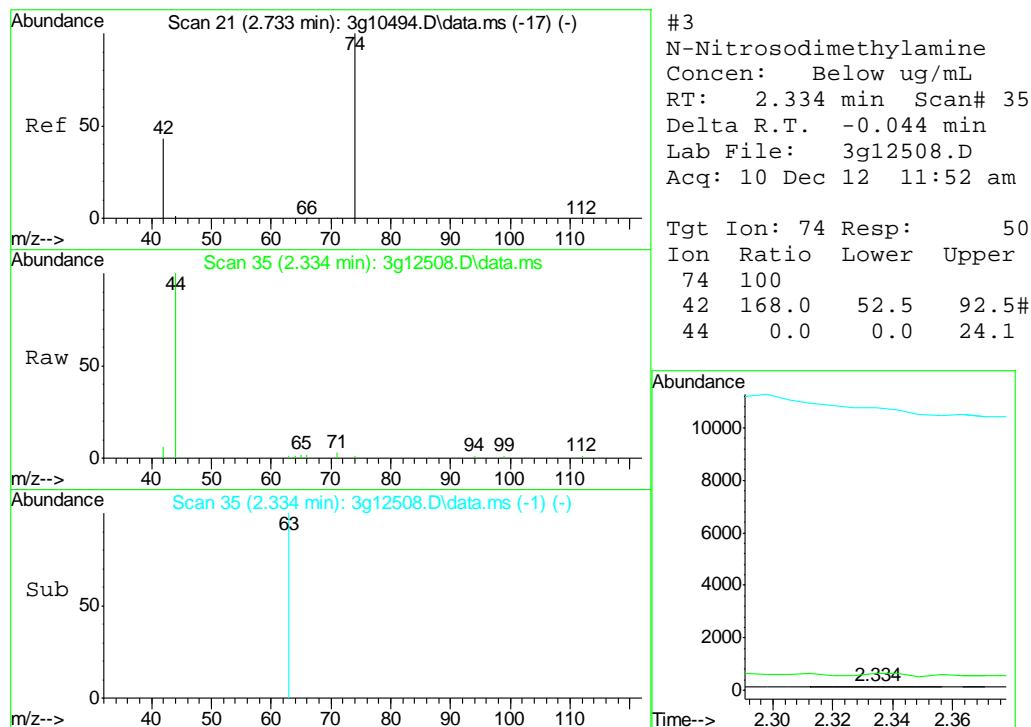
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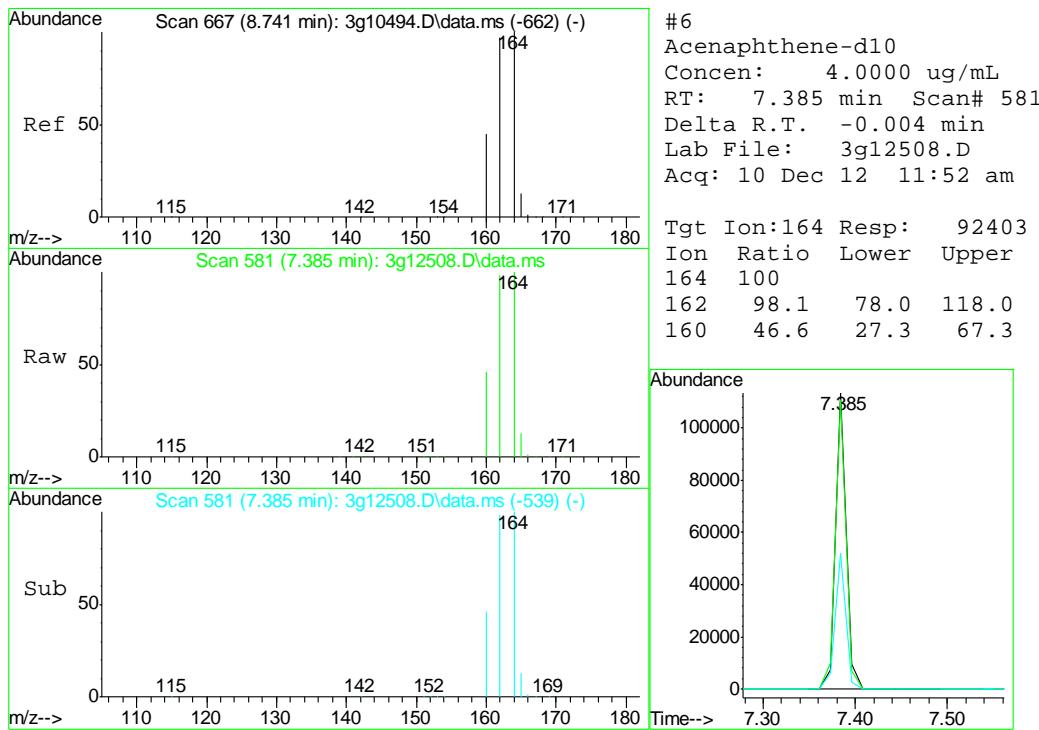
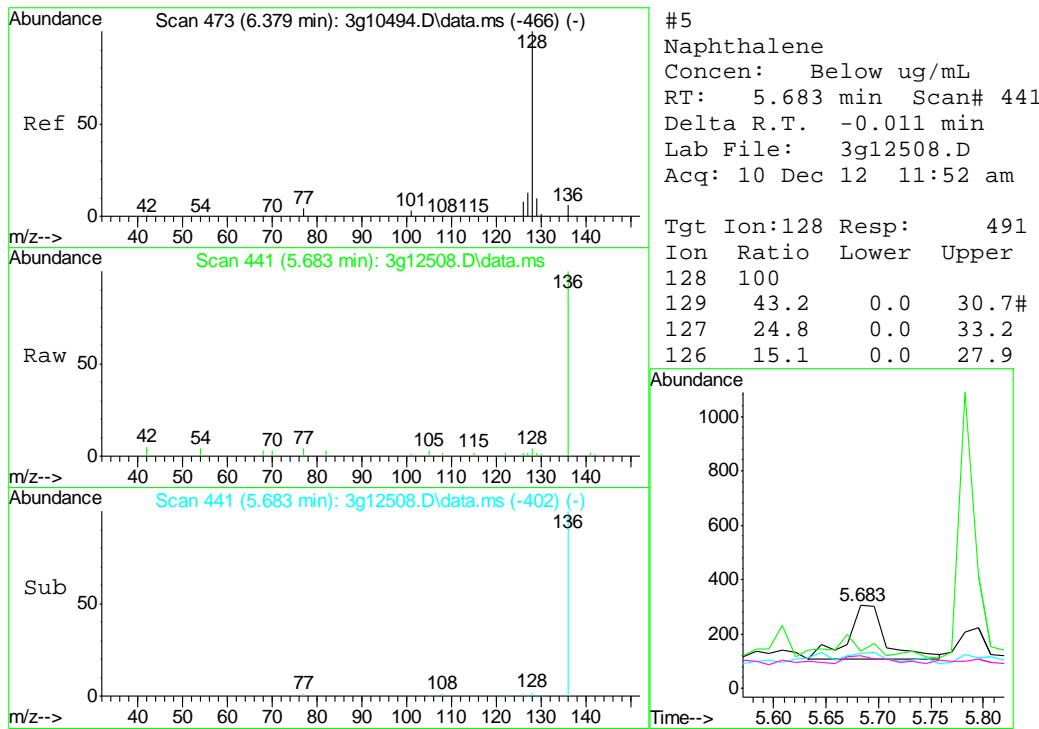
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 Misc : OP7075,E3G593,30.00,,,1,1  
 ALS Vial : 4 Sample Multiplier: 1

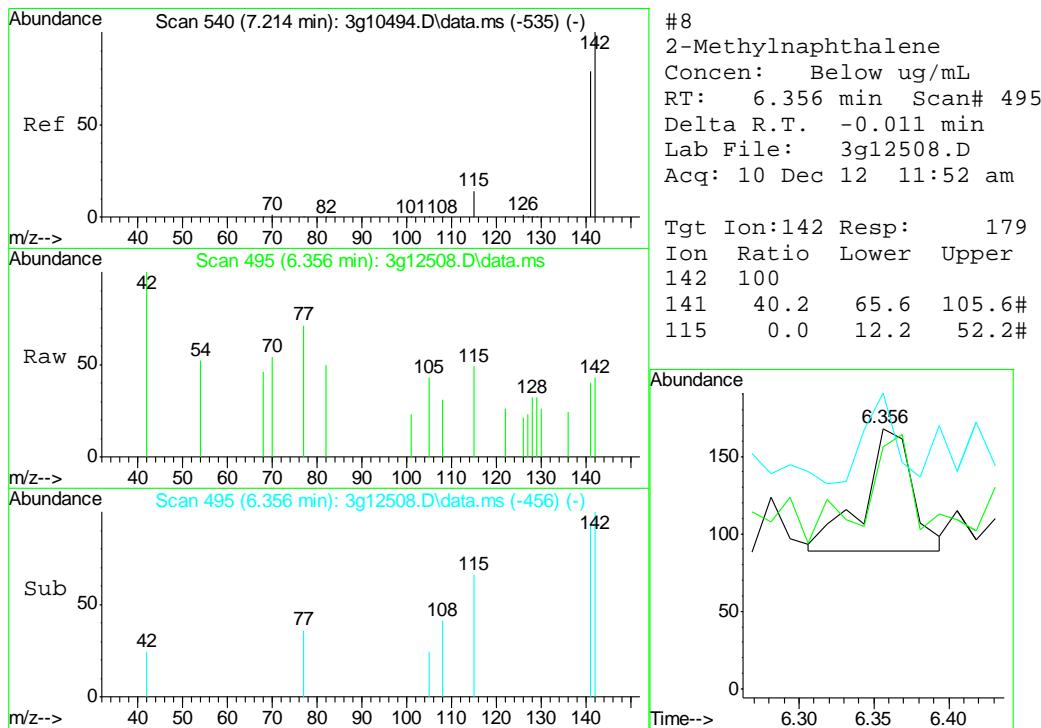
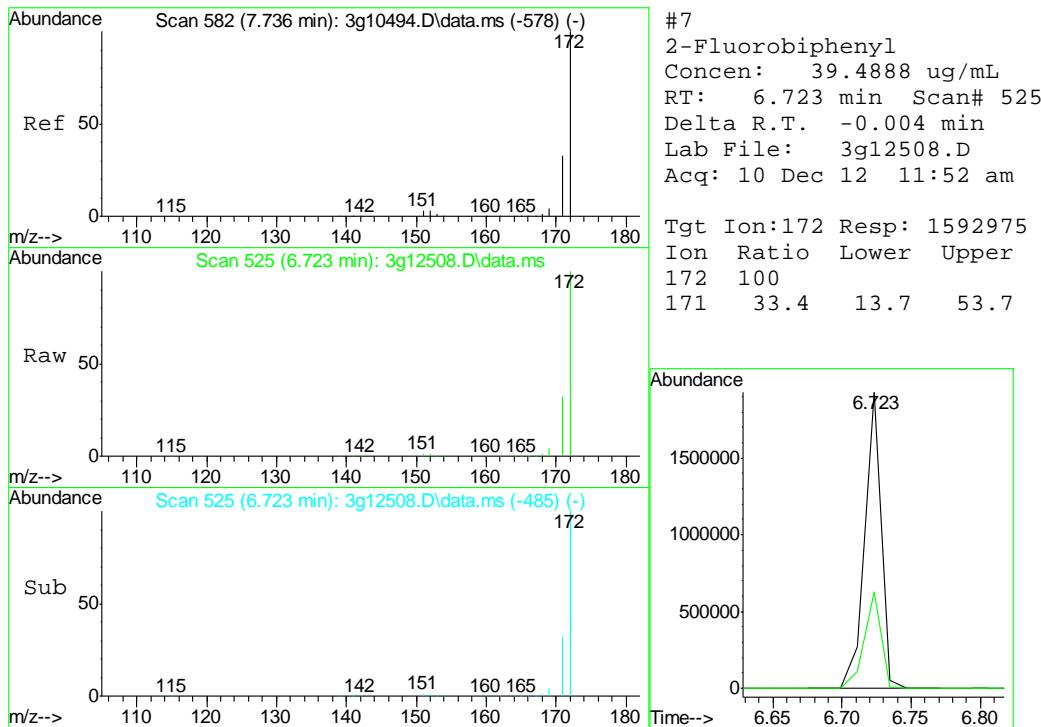
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 Response via : Initial Calibration

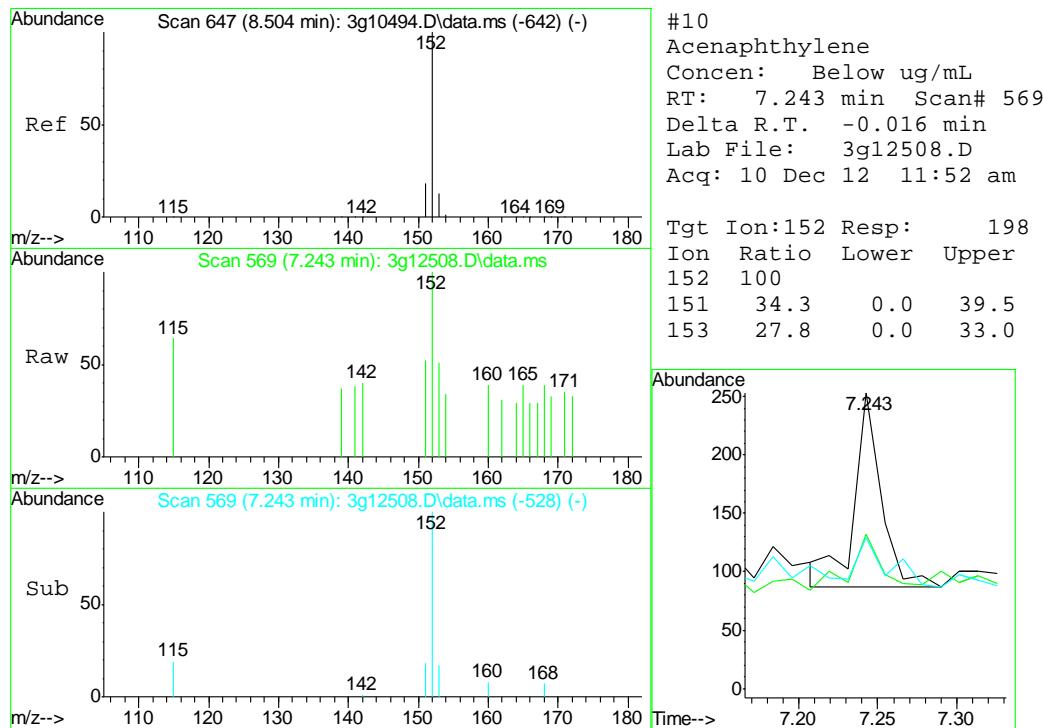
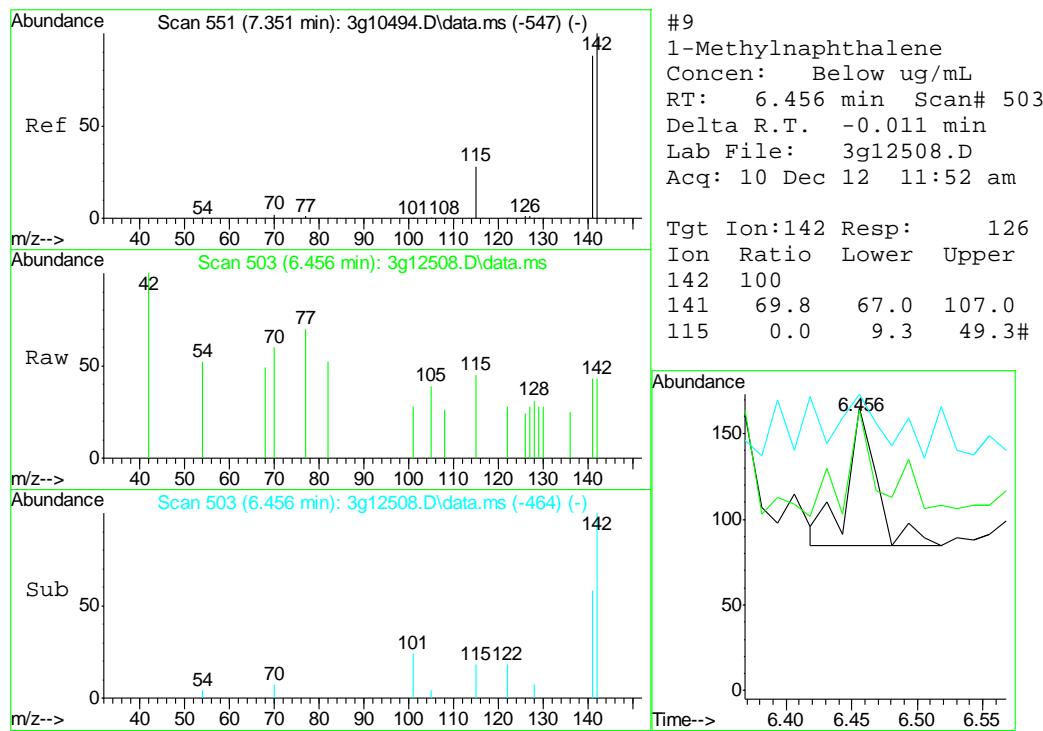


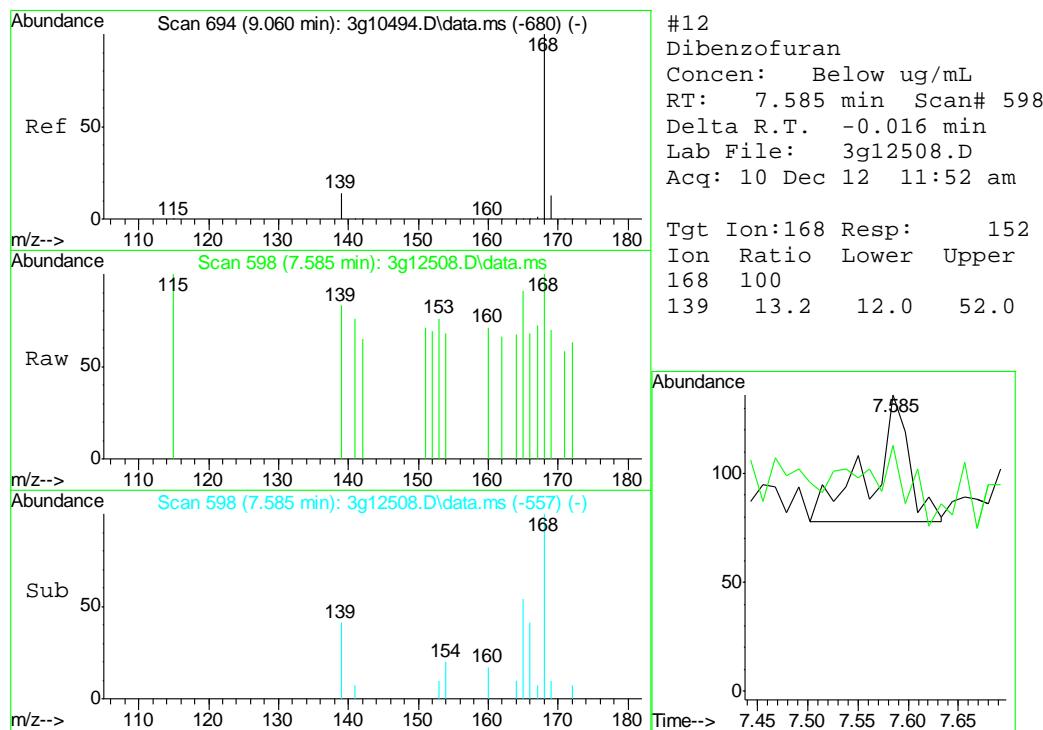
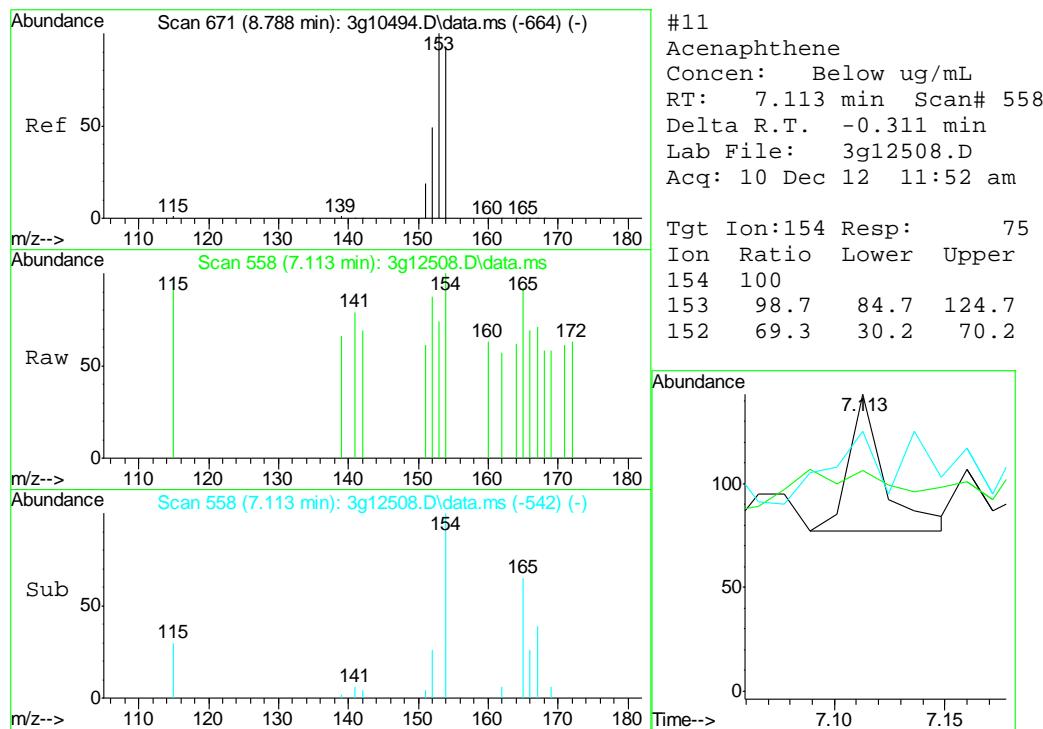


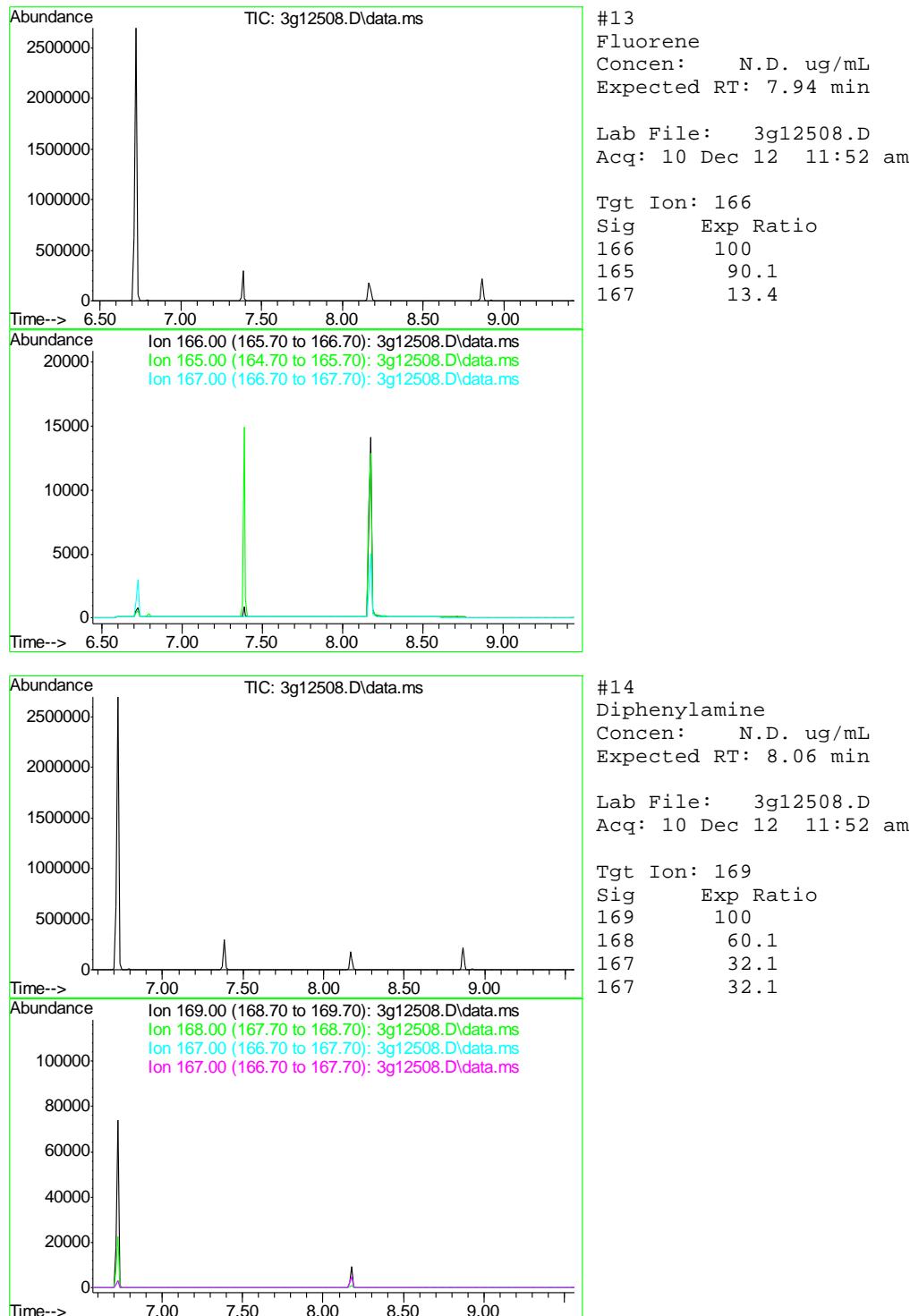


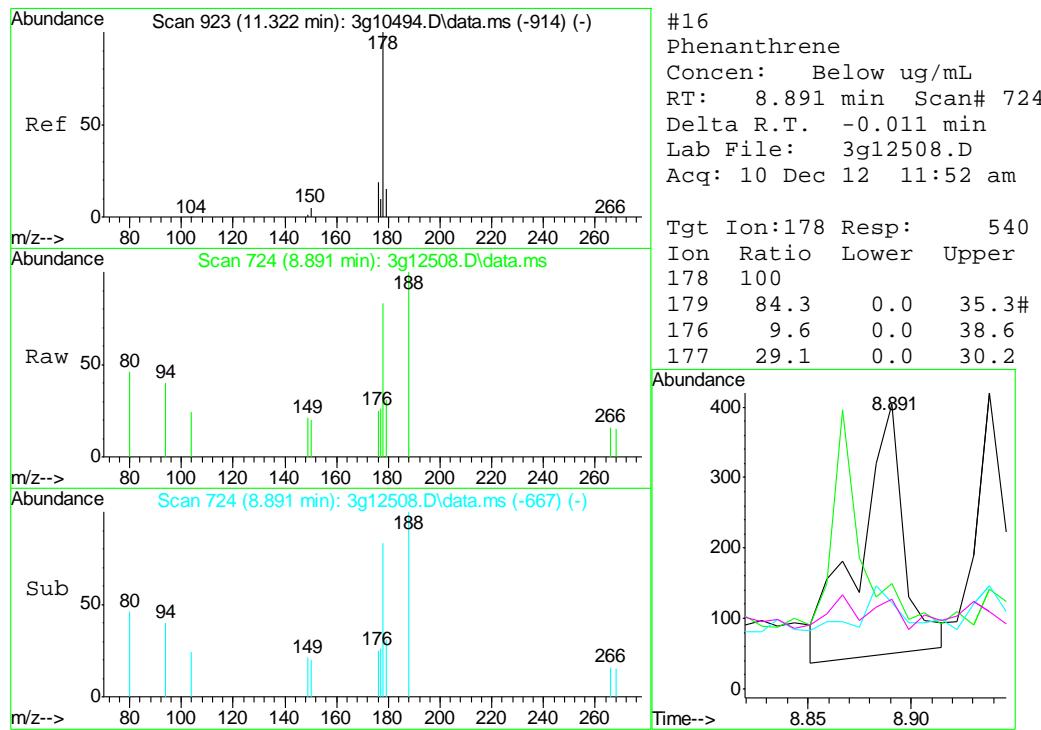
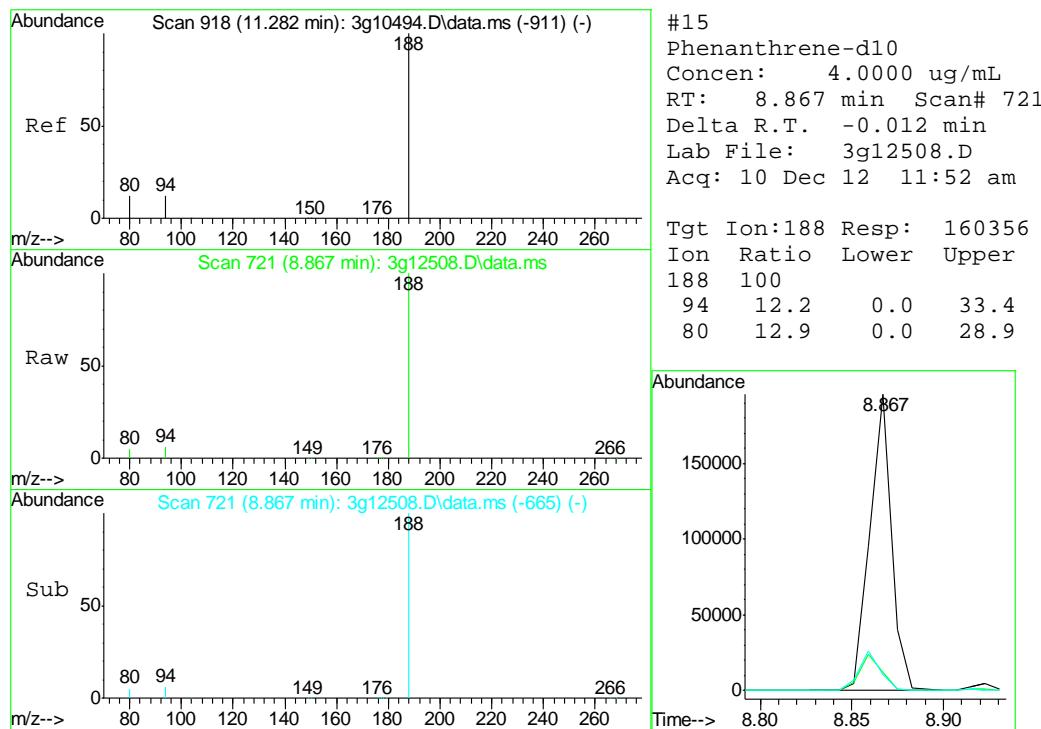


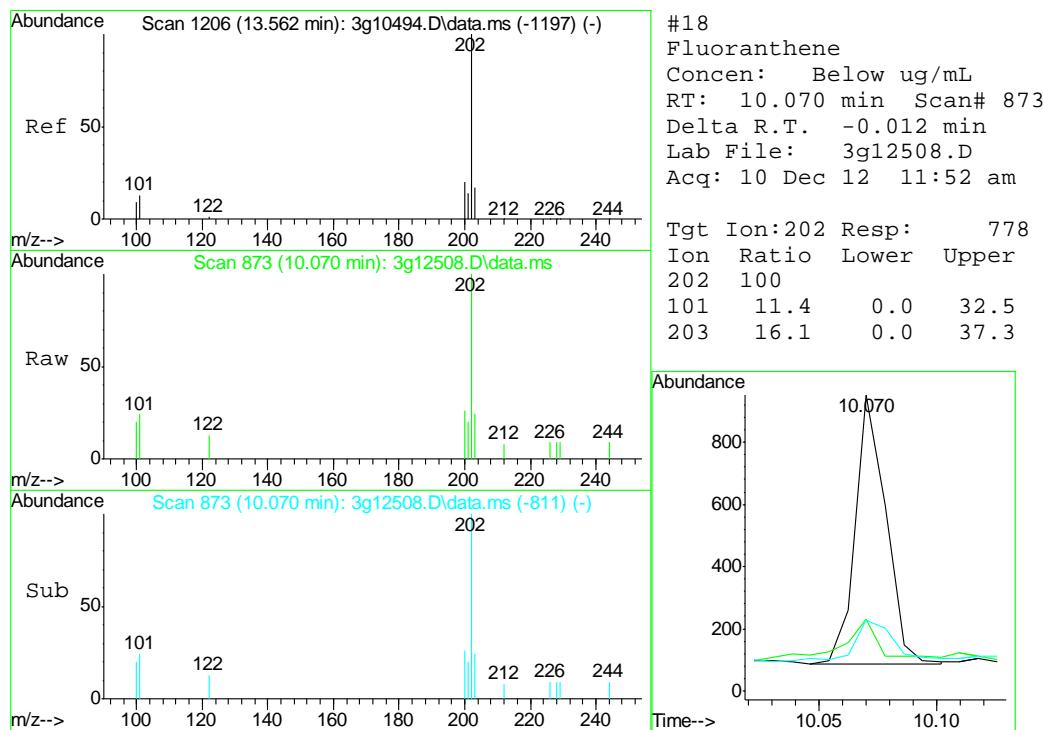
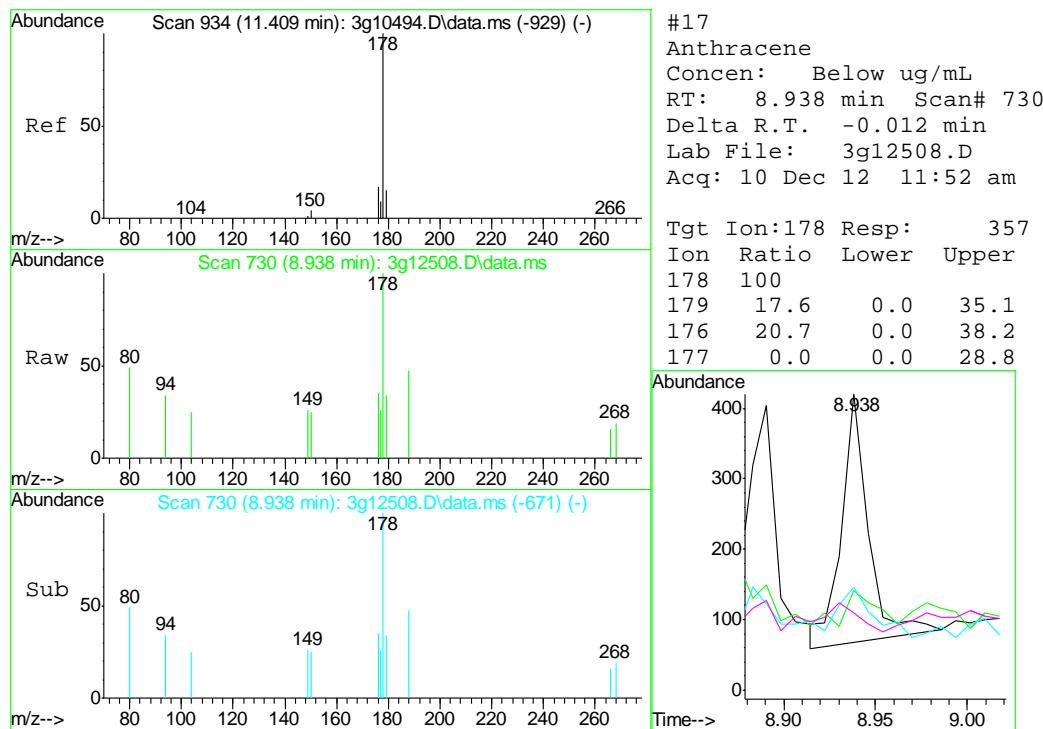


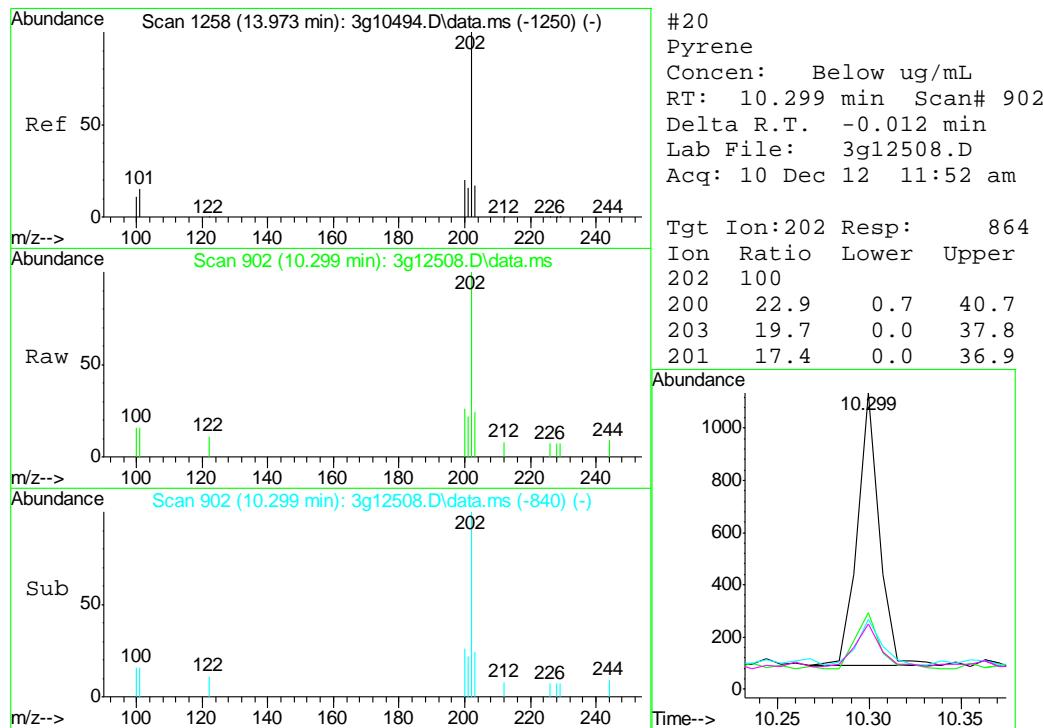
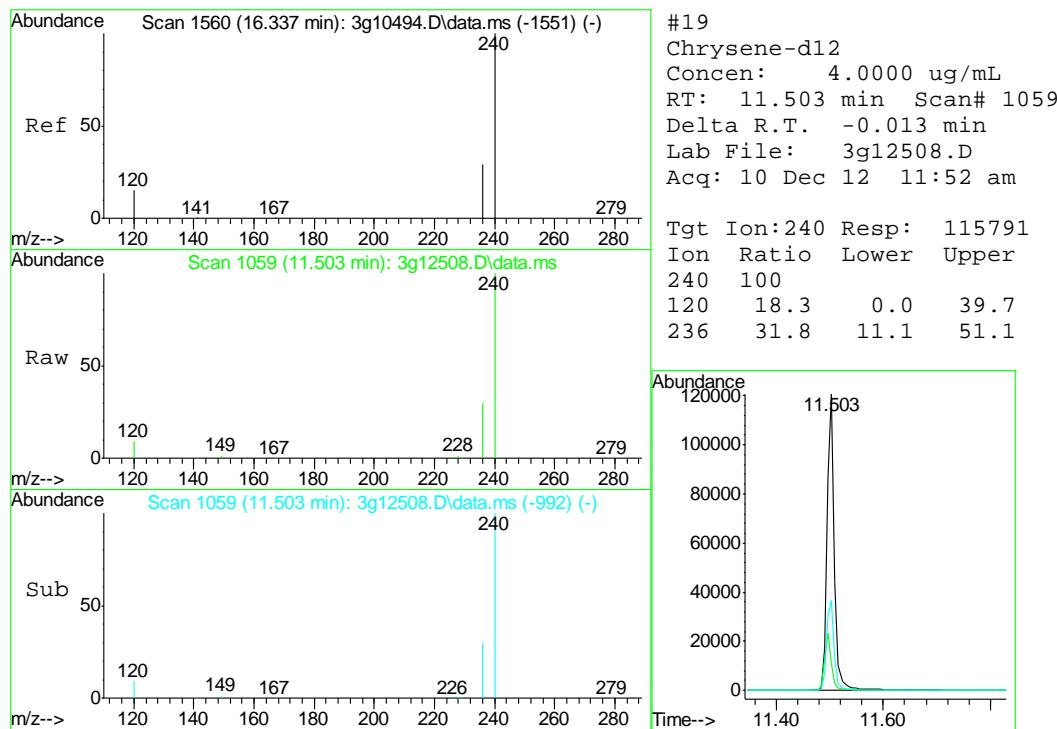


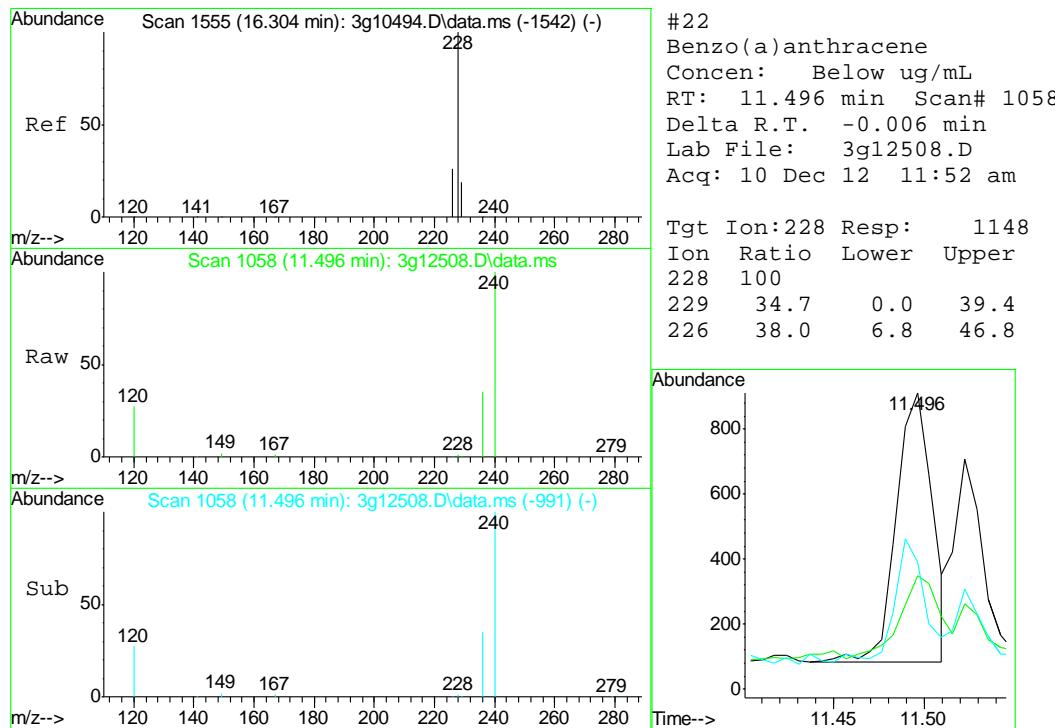
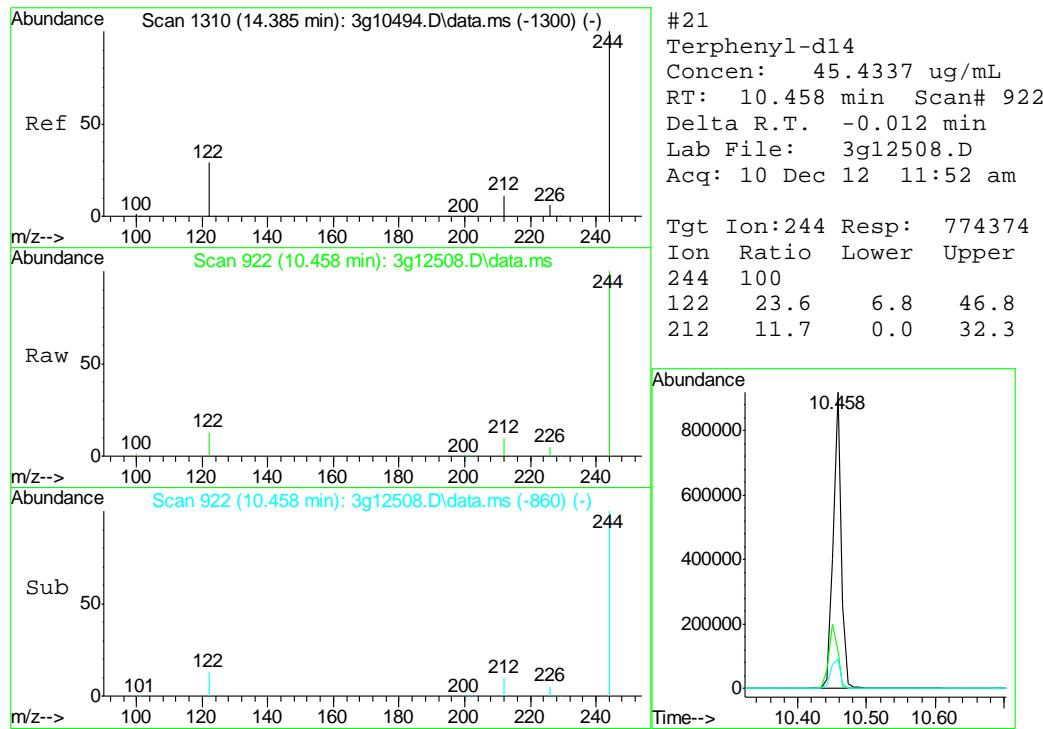


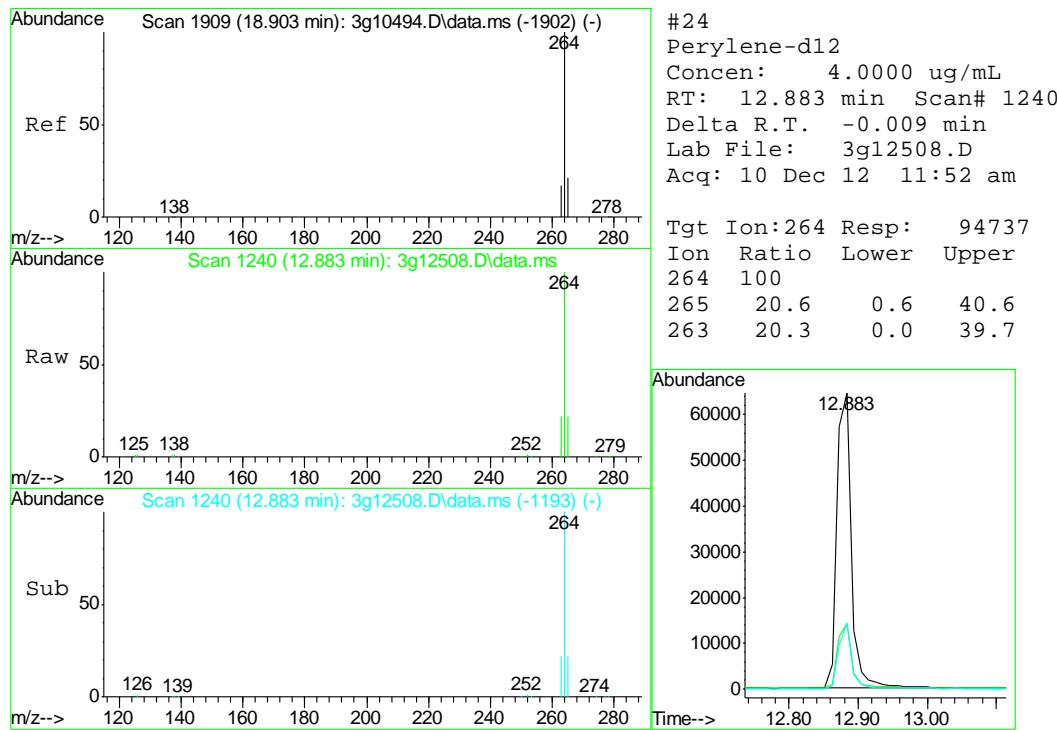
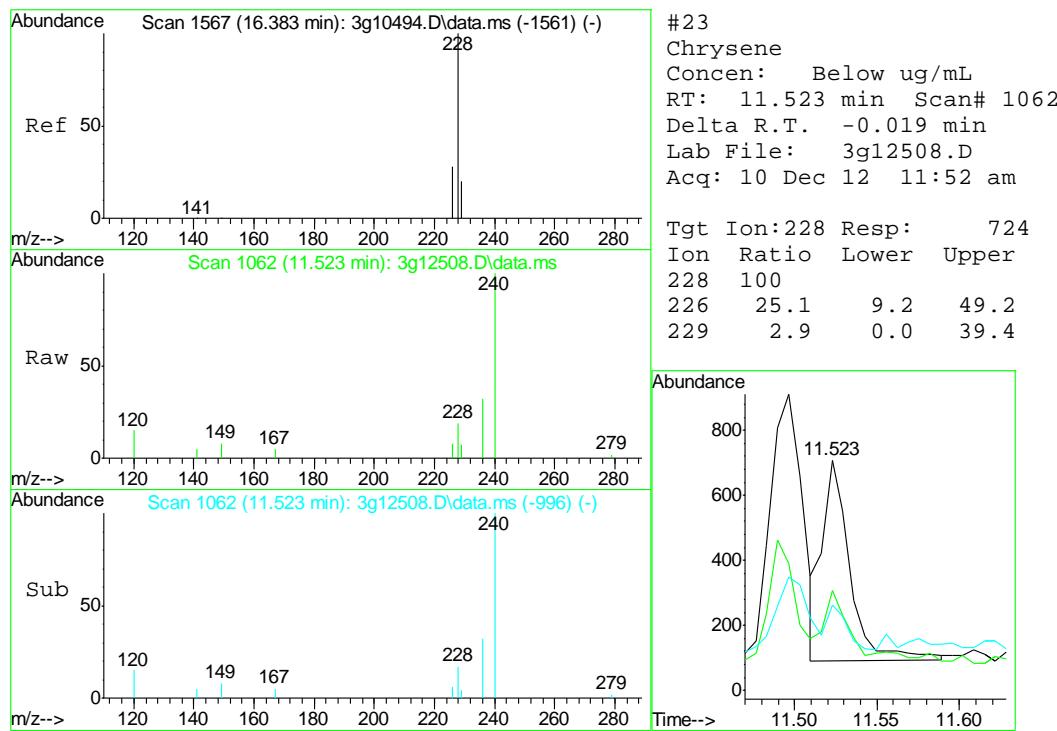


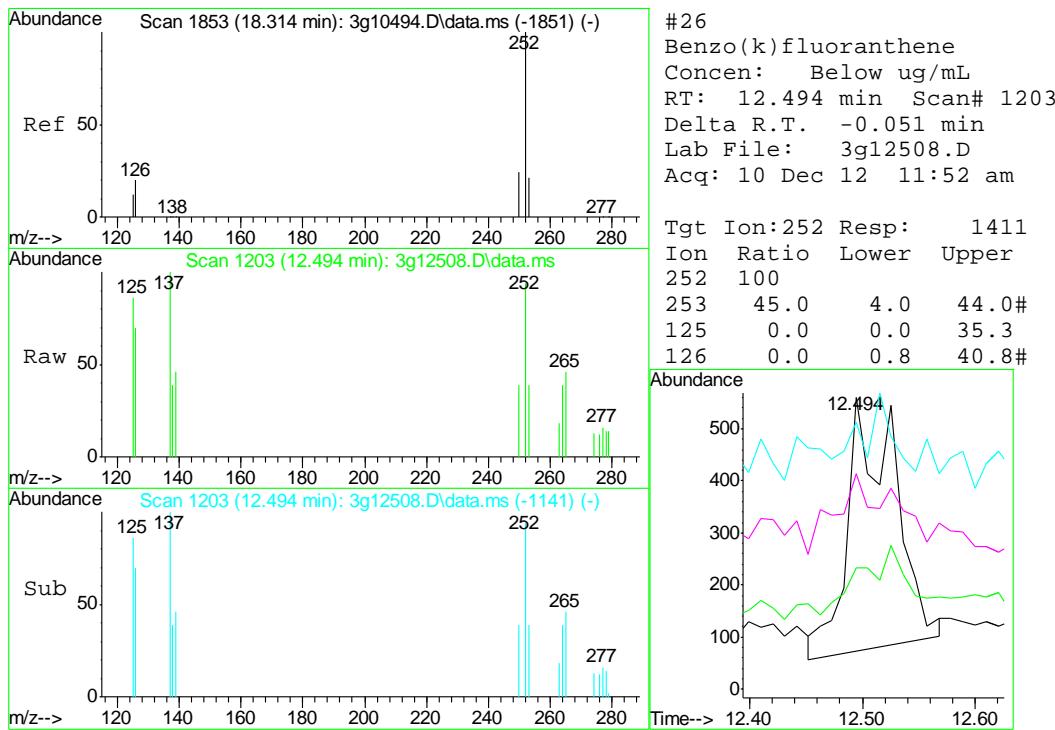
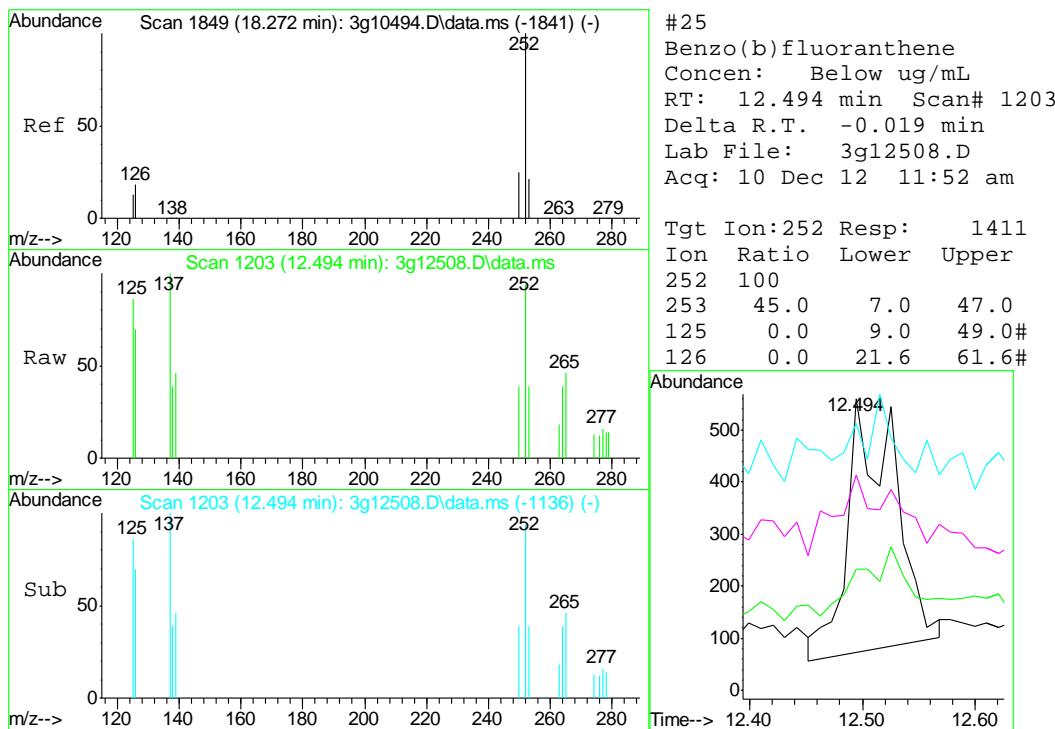


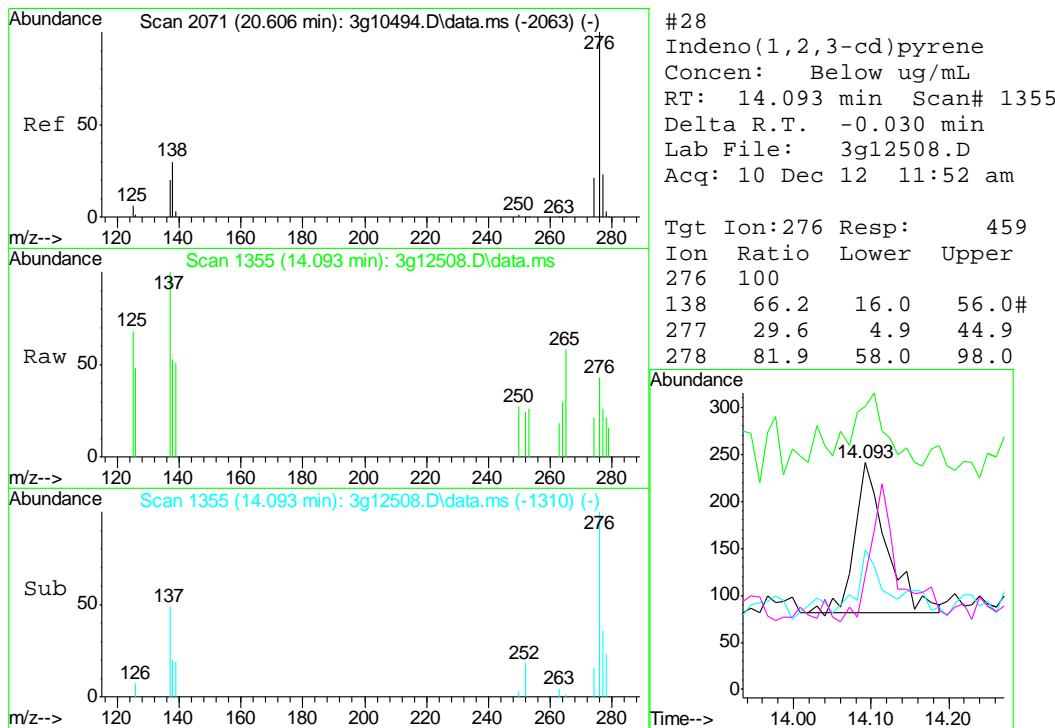
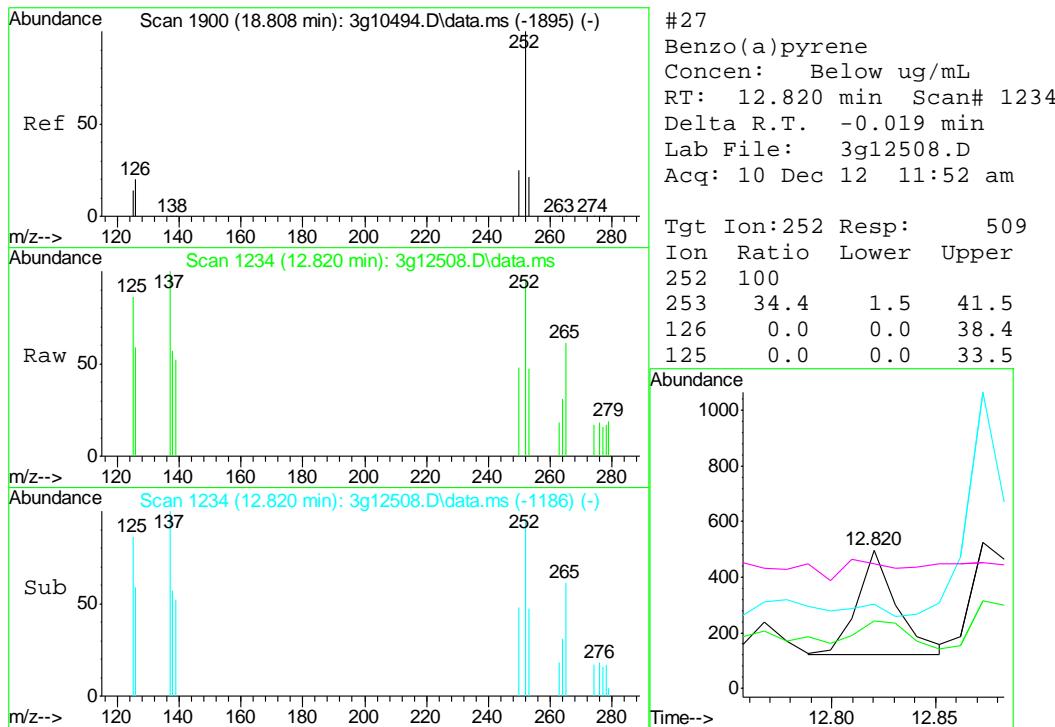


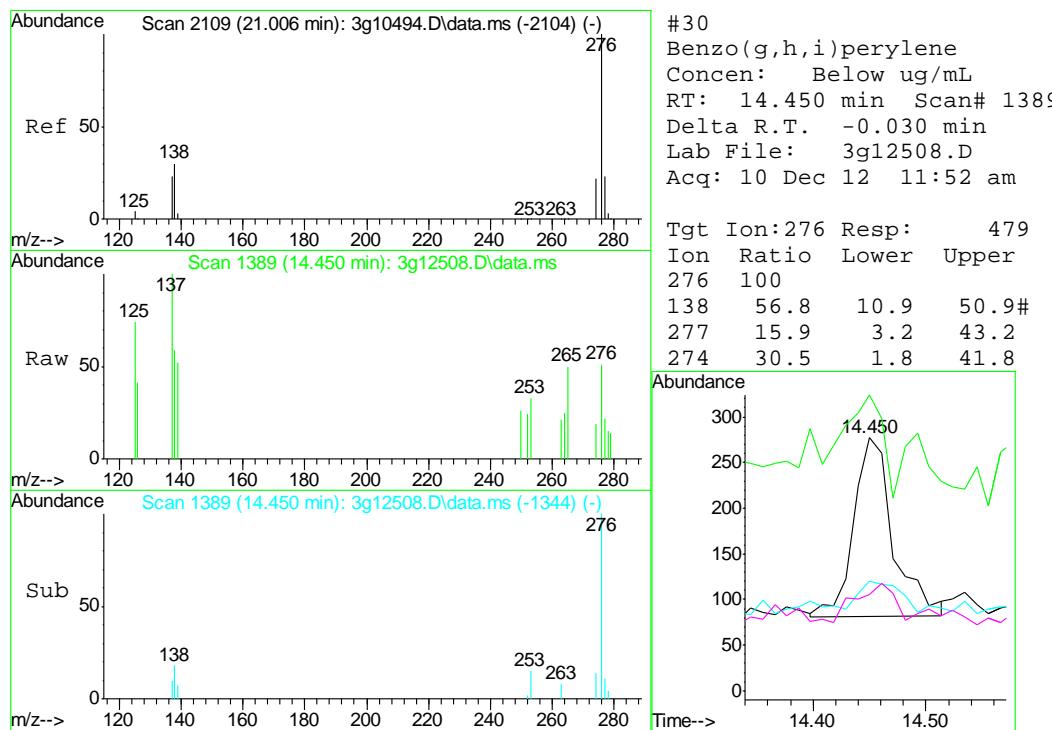
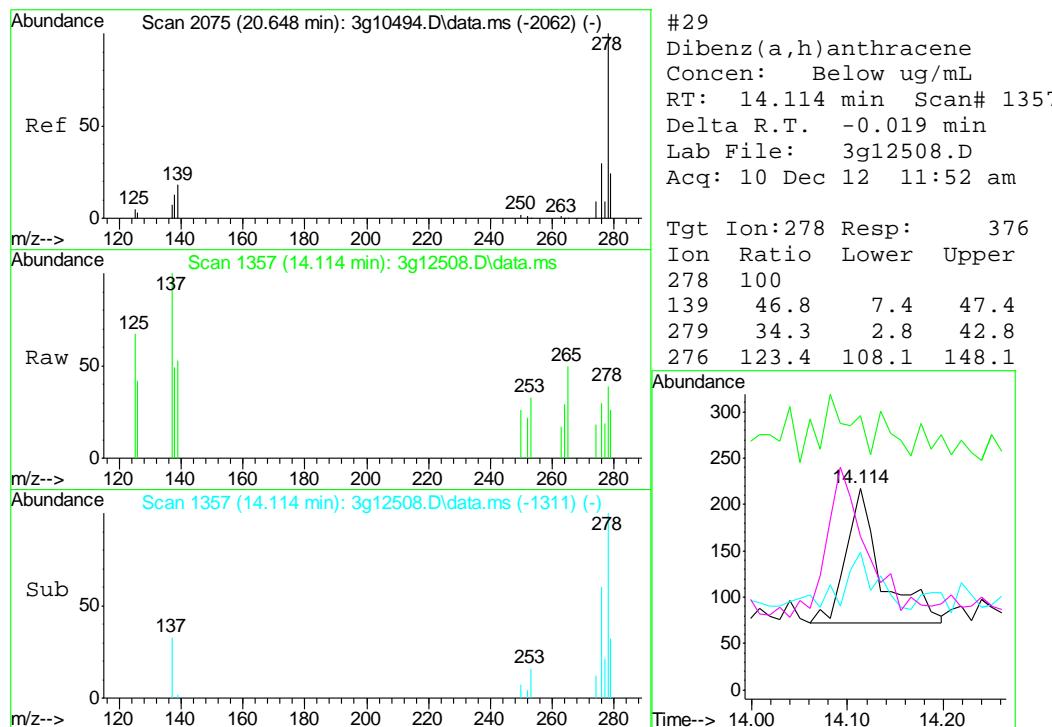














## GC Volatiles

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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:** D41506  
**Account:** XTOKWR XTO Energy  
**Project:** PCU 296-5A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
GGB1023-MB	GB18762.D	1	12/06/12	SK	n/a	n/a	GGB1023

The QC reported here applies to the following samples:

**Method:** SW846 8015B

D41506-1

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	89%      60-140%

10.1.1

10

## Blank Spike Summary

Page 1 of 1

Job Number: D41506

Account: XTOKWR XTO Energy

Project: PCU 296-5A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
GGB1023-BS	GB18763.D	1	12/06/12	SK	n/a	n/a	GGB1023

The QC reported here applies to the following samples:

Method: SW846 8015B

D41506-1

CAS No.	Compound	Spike mg/kg	BSP mg/kg	BSP %	Limits
	TPH-GRO (C6-C10)	110	123	112	70-130

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

10.2.1  
**10**

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\* = Outside of Control Limits.

# Matrix Spike/Matrix Spike Duplicate Summary

Page 1 of 1

Job Number: D41506

Account: XTOKWR XTO Energy

Project: PCU 296-5A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
D41506-1MS	GB18765.D	1	12/06/12	SK	n/a	n/a	GGB1023
D41506-1MSD	GB18766.D	1	12/06/12	SK	n/a	n/a	GGB1023
D41506-1	GB18764.D	1	12/06/12	SK	n/a	n/a	GGB1023

The QC reported here applies to the following samples:

Method: SW846 8015B

D41506-1

CAS No.	Compound	D41506-1		Spike	MS	MS	MSD	MSD	RPD	Limits Rec/RPD
		mg/kg	Q	mg/kg	mg/kg	%	mg/kg	%		
	TPH-GRO (C6-C10)	ND		132	147	111	148	112	1	70-130/30

CAS No.	Surrogate Recoveries	MS	MSD	D41506-1	Limits
120-82-1	1,2,4-Trichlorobenzene	96%	98%	90%	60-140%

\* = Outside of Control Limits.

10.3.1  
10



## GC Volatiles

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Raw Data

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## Quantitation Report (QT Reviewed)

Signal #1 : Y:\1\DATA\120612\GB18764.D\FID1A.CH Vial: 5  
 Signal #2 : Y:\1\DATA\120612\GB18764.D\FID2B.CH  
 Acq On : 6 Dec 2012 1:24 pm Operator: StephK  
 Sample : D41506-1, 50X Inst : GC/MS Ins  
 Misc : GC3278,GGB1023,,5.072,,100,5,1 Multiplr: 1.00  
 IntFile Signal #1: TVH1.E IntFile Signal #2: FB2.E  
 Quant Time: Dec 06 14:10:16 2012 Quant Results File: TB868GB868SOIL.RES

Quant Method : C:\MSDCHEM\1...\TB868GB868SOIL.M (Chemstation Integrator)  
 Title : 8015B/8021B TVH/BTEX  
 Last Update : Thu Dec 06 13:37:39 2012  
 Response via : Initial Calibration  
 DataAcq Meth : TVB4.M

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

Compound	R.T.	Response	Conc	Units
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System Monitoring Compounds

2) S	1,2,4-Trichlorobenzene	14.36	2825883	90.186 %
10) S	1,2,4-Trichlorobenzene (P)	14.36	14693615	90.407 %

Target Compounds

1) H	TVH-Gasoline	7.23	3723206	<MDL mg/L
4) T	Methyl-t-butyl-ether	0.00	0	N.D. ug/L d
5) T	Benzene	0.00	0	N.D. ug/L d
6) T	Toluene	7.66	221559	0.559 ug/L
7) T	Ethylbenzene	0.00	0	N.D. ug/L d
8) T	m,p-Xylene	10.47	175777	0.108 ug/L
9) T	o-Xylene	0.00	0	N.D. ug/L d
11) T	Naphthalene	14.54	423384	2.146 ug/L

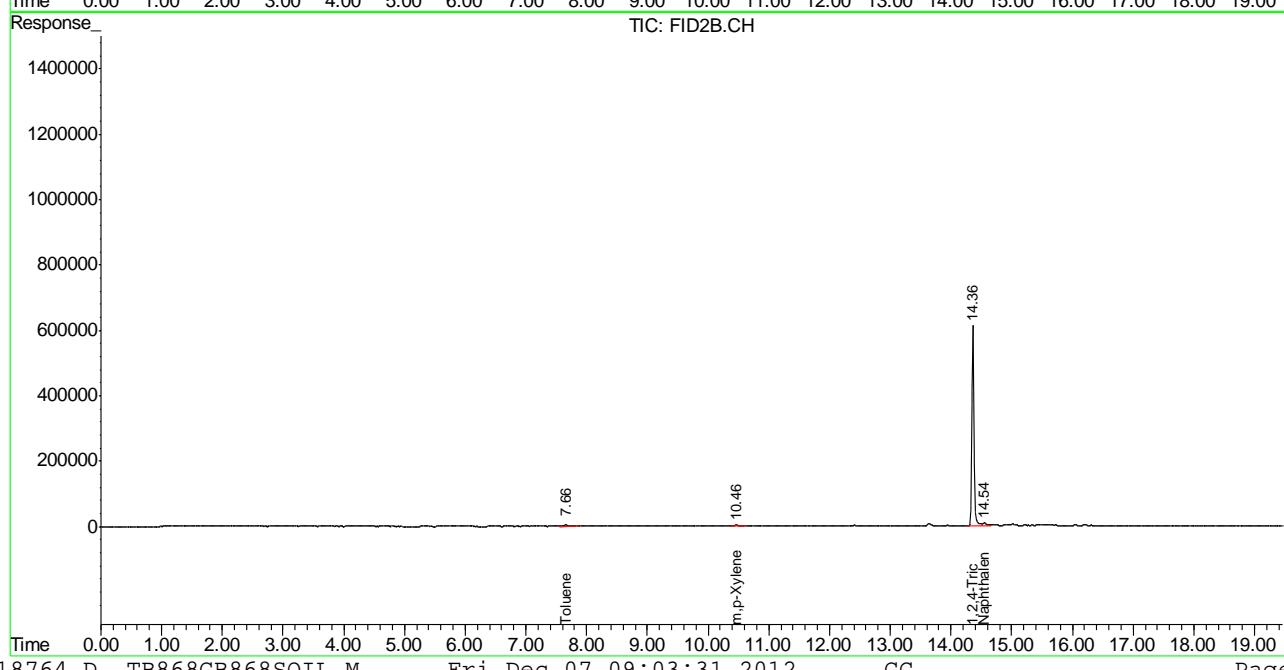
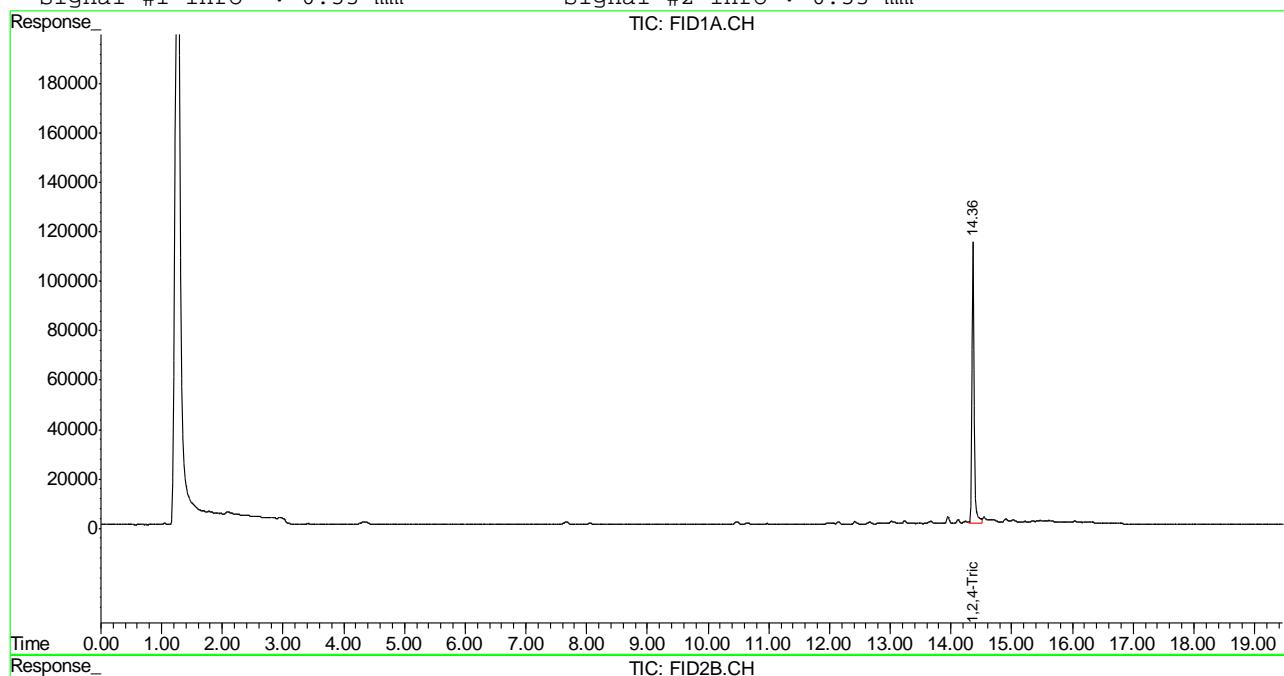
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 GB18764.D TB868GB868SOIL.M Fri Dec 07 09:03:31 2012 GC

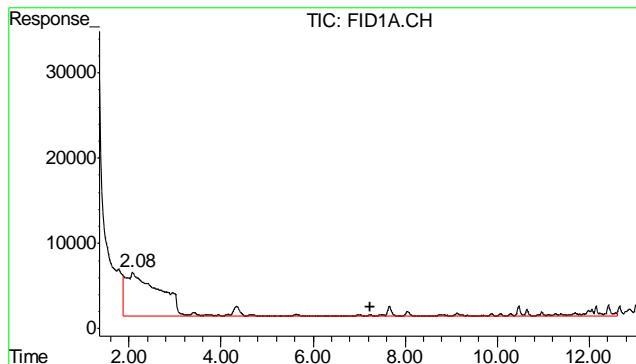
## Quantitation Report (QT Reviewed)

Signal #1 : Y:\1\DATA\120612\GB18764.D\FID1A.CH Vial: 5  
 Signal #2 : Y:\1\DATA\120612\GB18764.D\FID2B.CH  
 Acq On : 6 Dec 2012 1:24 pm Operator: StephK  
 Sample : D41506-1, 50X Inst : GC/MS Ins  
 Misc : GC3278,GGB1023,,5.072,,100,5,1 Multiplr: 1.00  
 IntFile Signal #1: TVH1.E IntFile Signal #2: FB2.E  
 Quant Time: Dec 6 14:10 2012 Quant Results File: TB868GB868SOIL.RES

Quant Method : C:\MSDCHEM\1...\TB868GB868SOIL.M (Chemstation Integrator)  
 Title : 8015B/8021B TVH/BTEX  
 Last Update : Thu Dec 06 13:37:39 2012  
 Response via : Multiple Level Calibration  
 DataAcq Meth : TVB4.M

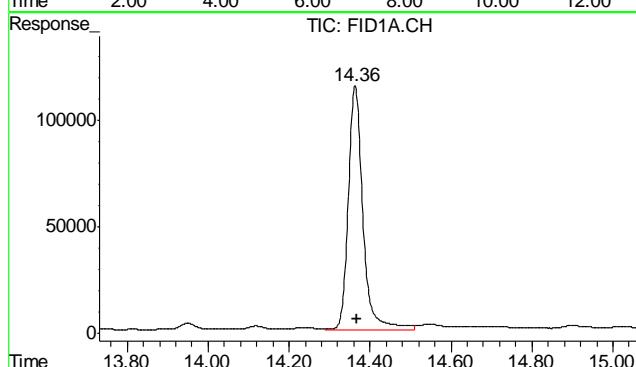
Volume Inj. :  
 Signal #1 Phase : DB-624 Signal #2 Phase: DB-624  
 Signal #1 Info : 0.53 mm Signal #2 Info : 0.53 mm





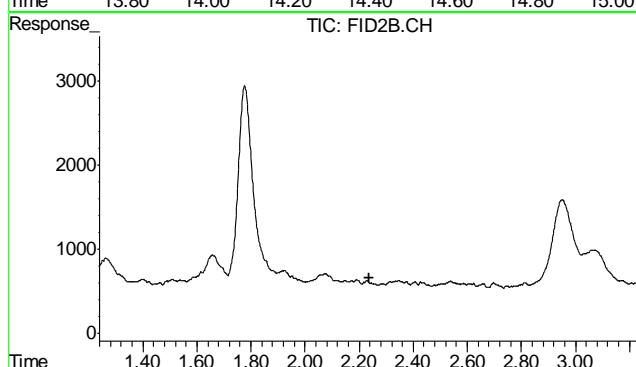
#1 TVH-Gasoline

R.T.: 7.230 min  
Delta R.T.: 0.000 min  
Response: 3723206  
Conc: N.D.



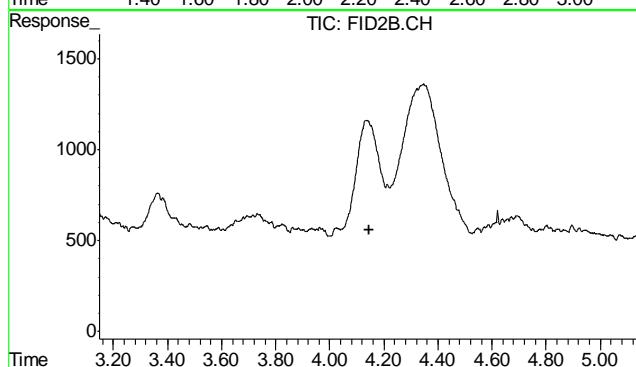
#2 1,2,4-Trichlorobenzene

R.T.: 14.363 min  
Delta R.T.: -0.004 min  
Response: 2825883  
Conc: 90.19 %



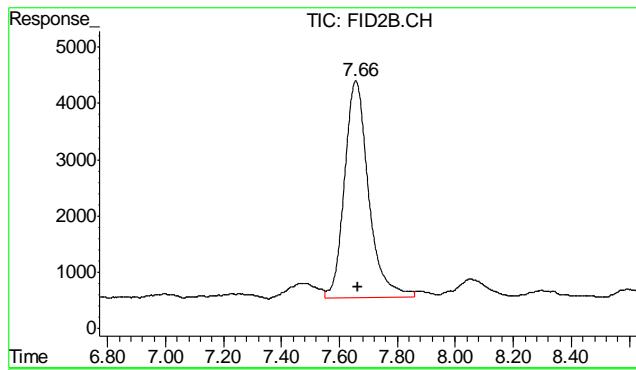
#4 Methyl-t-butyl-ether

R.T.: 0.000 min  
Exp R.T. : 2.239 min  
Response: 0  
Conc: N.D.



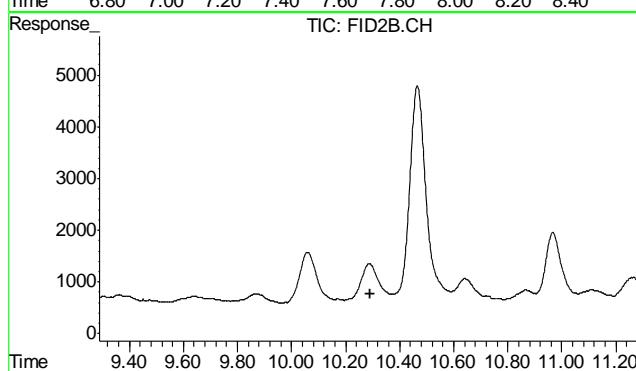
#5 Benzene

R.T.: 0.000 min  
Exp R.T. : 4.147 min  
Response: 0  
Conc: N.D.



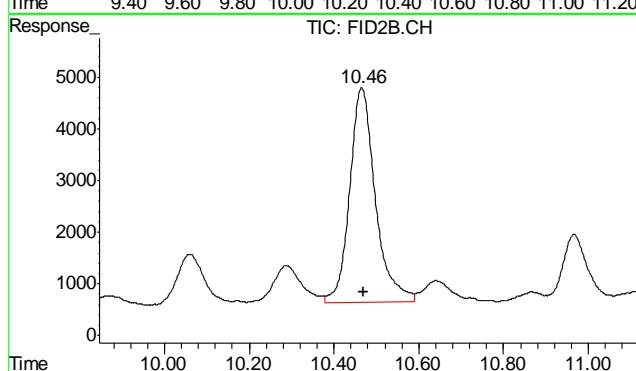
#6 Toluene

R.T.: 7.656 min  
 Delta R.T.: -0.008 min  
 Response: 221559  
 Conc: 0.56 ug/L



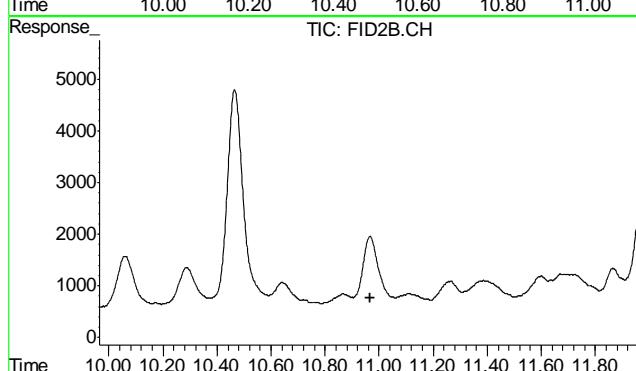
#7 Ethylbenzene

R.T.: 0.000 min  
 Exp R.T. : 10.290 min  
 Response: 0  
 Conc: N.D.



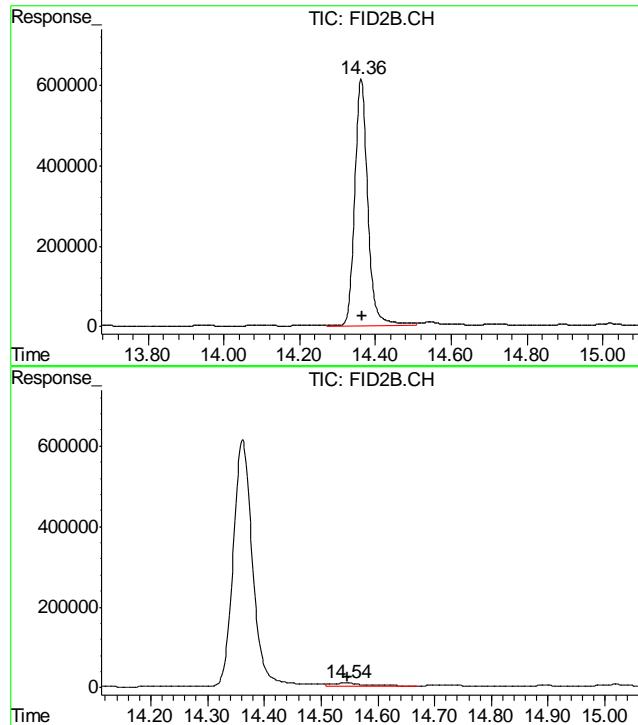
#8 m,p-Xylene

R.T.: 10.465 min  
 Delta R.T.: -0.004 min  
 Response: 175777  
 Conc: 0.11 ug/L



#9 o-Xylene

R.T.: 0.000 min  
 Exp R.T. : 10.965 min  
 Response: 0  
 Conc: N.D.



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

R.T.: 14.362 min  
Delta R.T.: -0.003 min  
Response: 14693615  
Conc: 90.41 %

#11 Naphthalene

R.T.: 14.544 min  
Delta R.T.: -0.003 min  
Response: 423384  
Conc: 2.15 ug/L

11.1

## Quantitation Report (QT Reviewed)

Signal #1 : Y:\1\DATA\120612\GB18762.D\FID1A.CH Vial: 3  
 Signal #2 : Y:\1\DATA\120612\GB18762.D\FID2B.CH  
 Acq On : 6 Dec 2012 12:14 pm Operator: StephK  
 Sample : MB Inst : GC/MS Ins  
 Misc : GC3278,GGB1023,5.000,,100,5,1 Multiplr: 1.00  
 IntFile Signal #1: TVH1.E IntFile Signal #2: FB2.E  
 Quant Time: Dec 06 13:37:59 2012 Quant Results File: TB868GB868SOIL.RES

Quant Method : C:\MSDCHEM\1...\TB868GB868SOIL.M (Chemstation Integrator)  
 Title : 8015B/8021B TVH/BTEX  
 Last Update : Thu Dec 06 13:37:39 2012  
 Response via : Initial Calibration  
 DataAcq Meth : TVB4.M

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

Compound	R.T.	Response	Conc	Units
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System Monitoring Compounds

2) S 1,2,4-Trichlorobenzene	14.37	2778216	88.665	%
10) S 1,2,4-Trichlorobenzene (P)	14.37	14665342	90.233	%

Target Compounds

1) H TVH-Gasoline	7.23	3435907	<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	7.67	165737	0.418	ug/L
7) T Ethylbenzene	0.00	0	N.D.	ug/L d
8) T m,p-Xylene	0.00	0	N.D.	ug/L d
9) T o-Xylene	0.00	0	N.D.	ug/L d
11) T Naphthalene	14.55	199919	1.013	ug/L

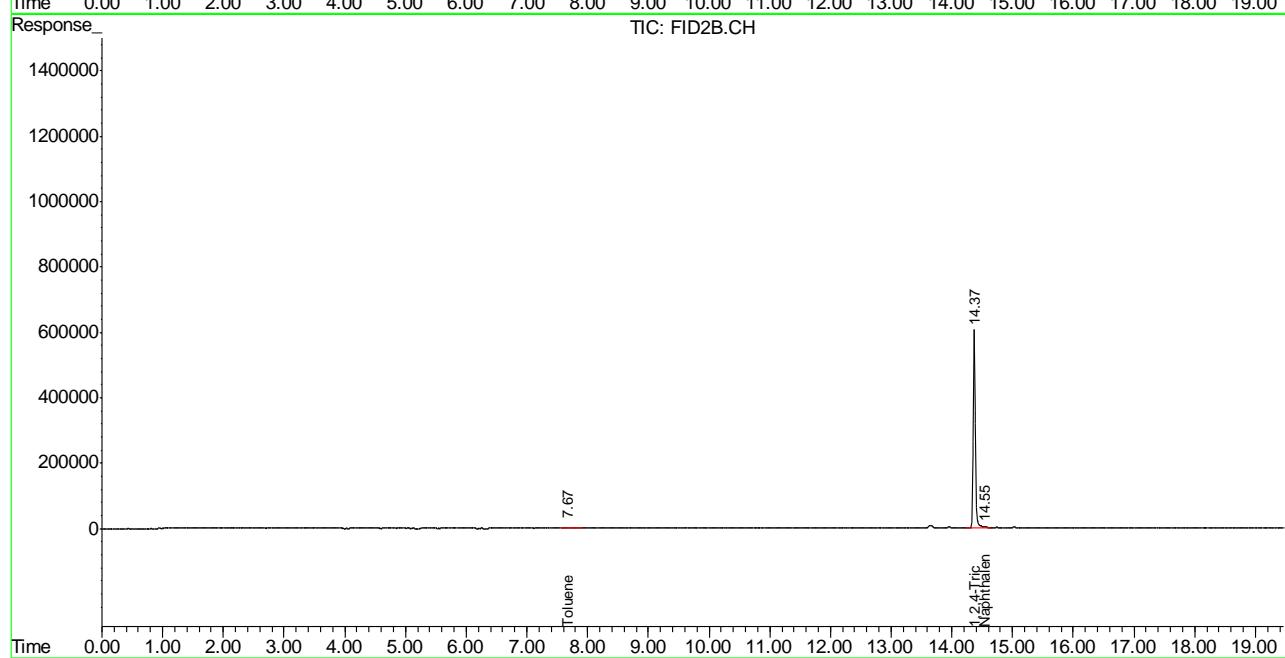
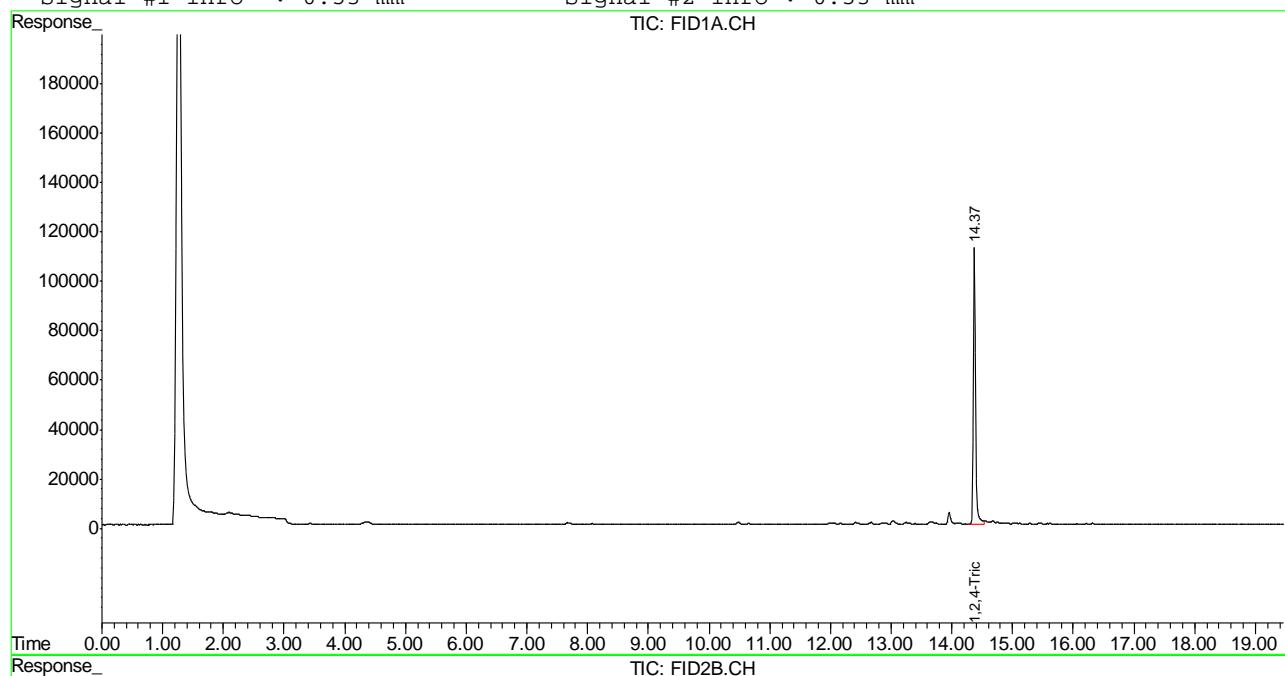
(f)=RT Delta > 1/2 Window (m)=manual int.  
 GB18762.D TB868GB868SOIL.M Fri Dec 07 09:03:25 2012 GC

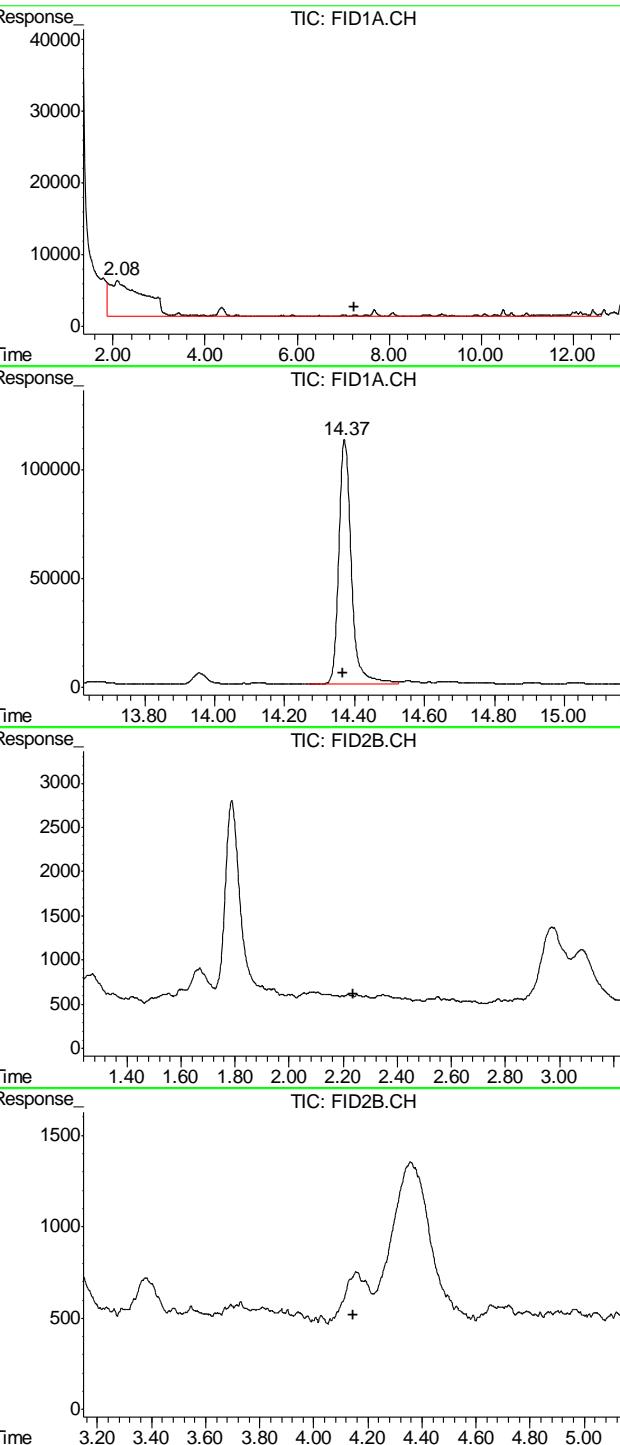
## Quantitation Report (QT Reviewed)

Signal #1 : Y:\1\DATA\120612\GB18762.D\FID1A.CH Vial: 3  
 Signal #2 : Y:\1\DATA\120612\GB18762.D\FID2B.CH  
 Acq On : 6 Dec 2012 12:14 pm Operator: StephK  
 Sample : MB Inst : GC/MS Ins  
 Misc : GC3278,GGB1023,5.000,,100,5,1 Multiplr: 1.00  
 IntFile Signal #1: TVH1.E IntFile Signal #2: FB2.E  
 Quant Time: Dec 6 13:38 2012 Quant Results File: TB868GB868SOIL.RES

Quant Method : C:\MSDCHEM\1...\TB868GB868SOIL.M (Chemstation Integrator)  
 Title : 8015B/8021B TVH/BTEX  
 Last Update : Thu Dec 06 13:37:39 2012  
 Response via : Multiple Level Calibration  
 DataAcq Meth : TVB4.M

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



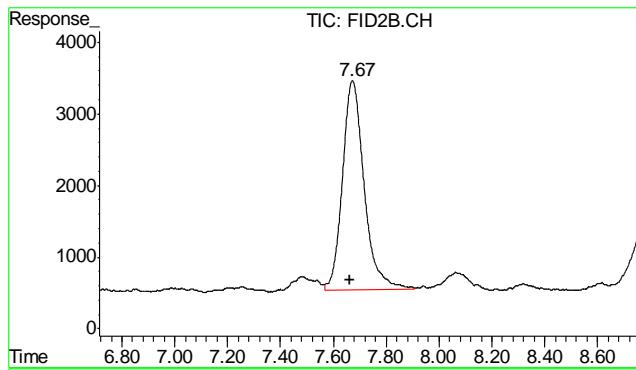


#1 TVH-Gasoline  
 R.T.: 7.230 min  
 Delta R.T.: 0.000 min  
 Response: 3435907  
 Conc: N.D.

#2 1,2,4-Trichlorobenzene  
 R.T.: 14.372 min  
 Delta R.T.: 0.004 min  
 Response: 2778216  
 Conc: 88.66 %

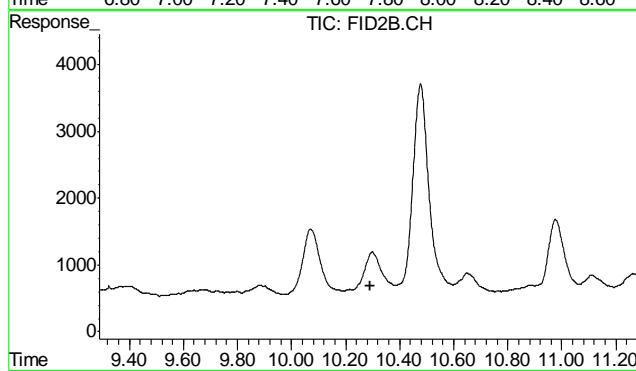
#4 Methyl-t-butyl-ether  
 R.T.: 0.000 min  
 Exp R.T. : 2.239 min  
 Response: 0  
 Conc: N.D.

#5 Benzene  
 R.T.: 0.000 min  
 Exp R.T. : 4.147 min  
 Response: 0  
 Conc: N.D.



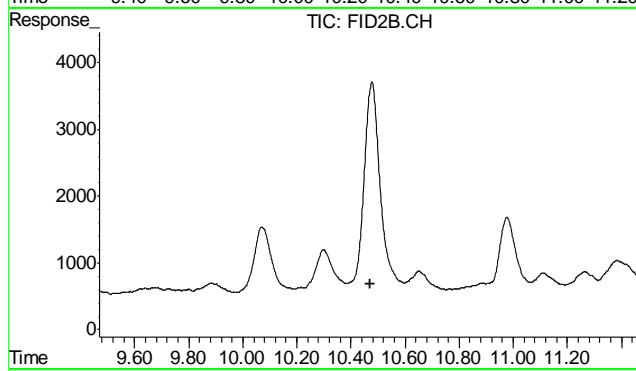
#6 Toluene

R.T.: 7.672 min  
Delta R.T.: 0.008 min  
Response: 165737  
Conc: 0.42 ug/L



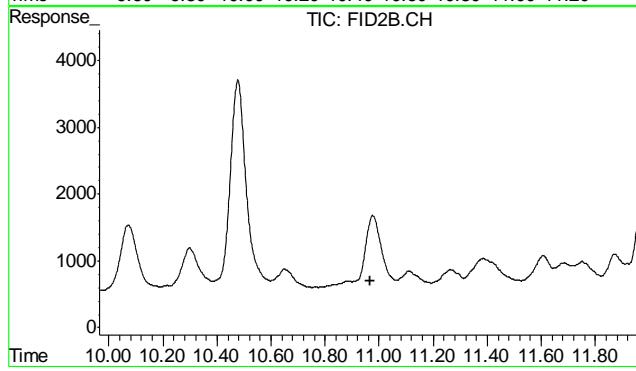
#7 Ethylbenzene

R.T.: 0.000 min  
Exp R.T. : 10.290 min  
Response: 0  
Conc: N.D.



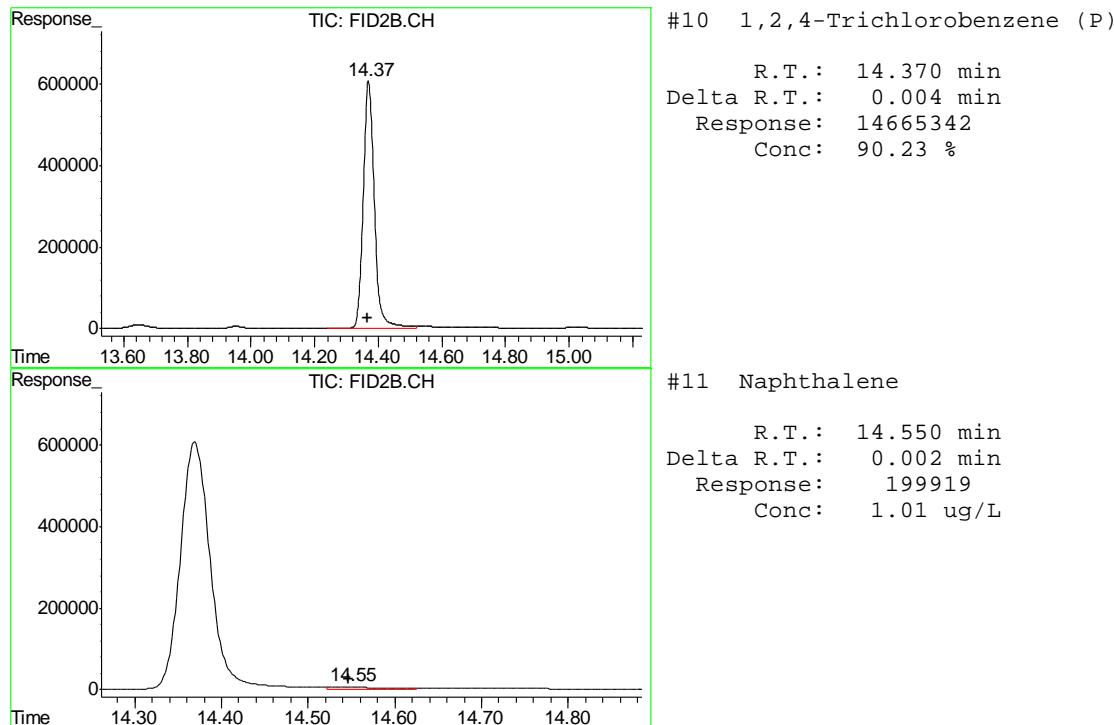
#8 m,p-Xylene

R.T.: 0.000 min  
Exp R.T. : 10.469 min  
Response: 0  
Conc: N.D.



#9 o-Xylene

R.T.: 0.000 min  
Exp R.T. : 10.965 min  
Response: 0  
Conc: N.D.





## 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:** D41506  
**Account:** XTOKWR XTO Energy  
**Project:** PCU 296-5A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
OP7061-MB	FD20235.D	1	12/07/12	AV	12/06/12	OP7061	GFD1015

The QC reported here applies to the following samples:

**Method:** SW846-8015B

D41506-1

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

CAS No.	Surrogate Recoveries	Limits
84-15-1	o-Terphenyl	78% 35-130%

## Blank Spike Summary

Page 1 of 1

Job Number: D41506

Account: XTOKWR XTO Energy

Project: PCU 296-5A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
OP7061-BS	FD20237.D	1	12/07/12	AV	12/06/12	OP7061	GFD1015

The QC reported here applies to the following samples:

Method: SW846-8015B

D41506-1

CAS No.	Compound	Spike mg/kg	BSP mg/kg	BSP %	Limits
	TPH-DRO (C10-C28)	667	695	104	48-130

CAS No.	Surrogate Recoveries	BSP	Limits
84-15-1	o-Terphenyl	99%	35-130%

\* = Outside of Control Limits.

12.2.1

12

# Matrix Spike/Matrix Spike Duplicate Summary

Page 1 of 1

Job Number: D41506

Account: XTOKWR XTO Energy

Project: PCU 296-5A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
OP7061-MS	FD20239.D	1	12/07/12	AV	12/06/12	OP7061	GFD1015
OP7061-MSD	FD20241.D	1	12/07/12	AV	12/06/12	OP7061	GFD1015
D41506-1	FD20243.D	1	12/07/12	AV	12/06/12	OP7061	GFD1015

The QC reported here applies to the following samples:

Method: SW846-8015B

D41506-1

CAS No.	Compound	D41506-1		Spike	MS	MS	MSD	MSD	RPD	Limits Rec/RPD
		mg/kg	Q	mg/kg	mg/kg	%	mg/kg	%		
	TPH-DRO (C10-C28)	33.8		739	699	90	732	94	5	20-168/30

CAS No.	Surrogate Recoveries	MS	MSD	D41506-1	Limits
84-15-1	o-Terphenyl	91%	89%	68%	35-130%

\* = Outside of Control Limits.

12.3.1  
12



## GC Semi-volatiles

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Raw Data

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## Quantitation Report (QT Reviewed)

Data File : C:\MSDCHEM\2\DATA\2012\DEC\FD120712\FD20243.D Vial: 7  
 Acq On : 07 Dec 2012 11:12 am Operator: ashleyv  
 Sample : D41506-1 Inst : FID5  
 Misc : OP7061,GFD1015,30.00,,,1,1 Multiplr: 1.00  
 IntFile : autoint1.e  
 Quant Time: Dec 07 11:37:25 2012 Quant Results File: DRO-GFD982F.RES

Quant Method : C:\MSDCHEM\2...\DRO-GFD982F.M (Chemstation Integrator)  
 Title : 8015B TEH  
 Last Update : Fri Nov 16 10:24:56 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
<hr/>				
System Monitoring Compounds				
1) S O-Terphenyl	8.93	75762260	1369.270	mg/L
<hr/>				
Target Compounds				
2) H TPH-DRO (c10-c28)	6.89	34755383	915.201	mg/L

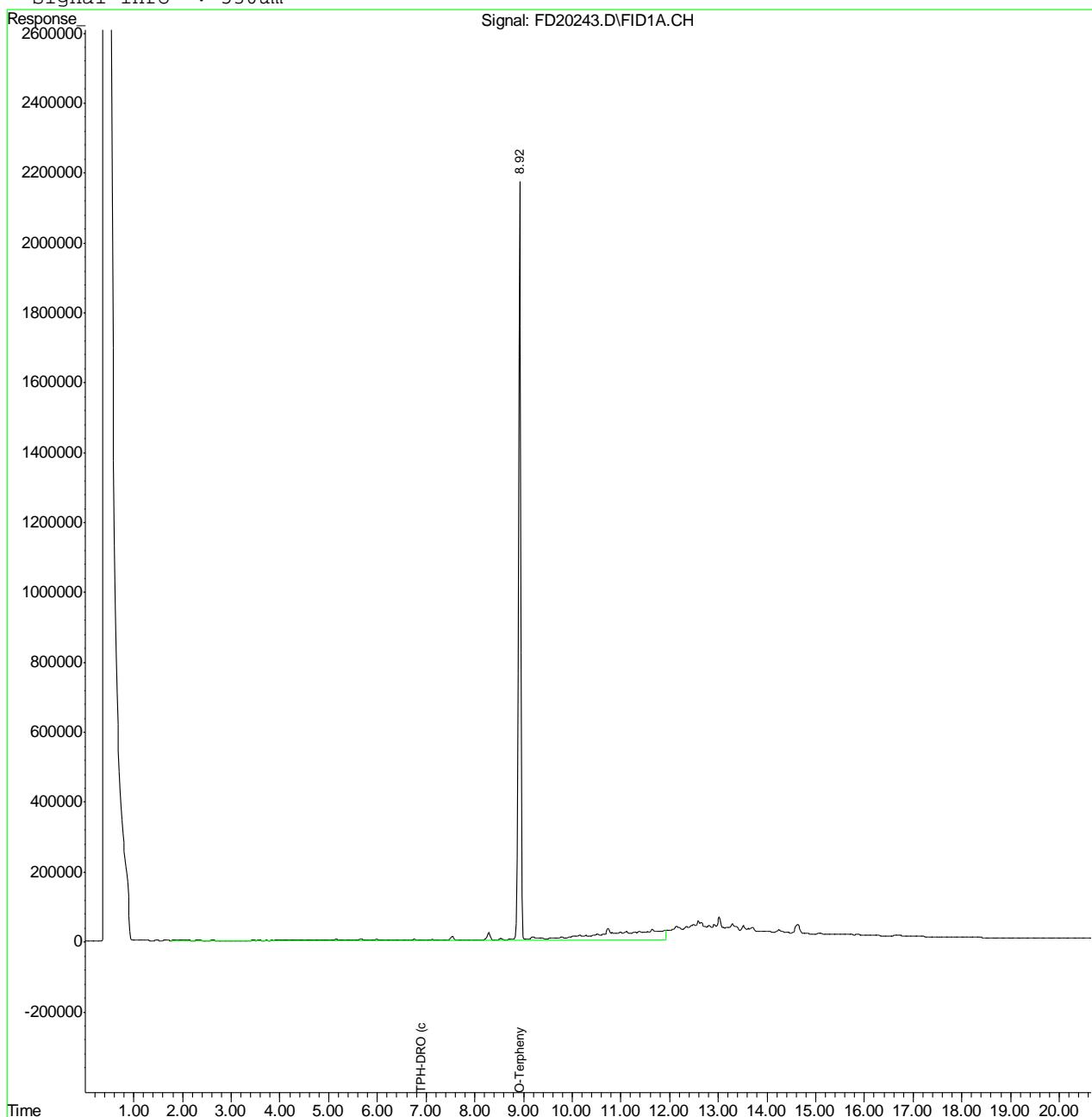
(f)=RT Delta > 1/2 Window (m)=manual int.  
 FD20243.D DRO-GFD982F.M Fri Dec 07 12:23:50 2012 GC

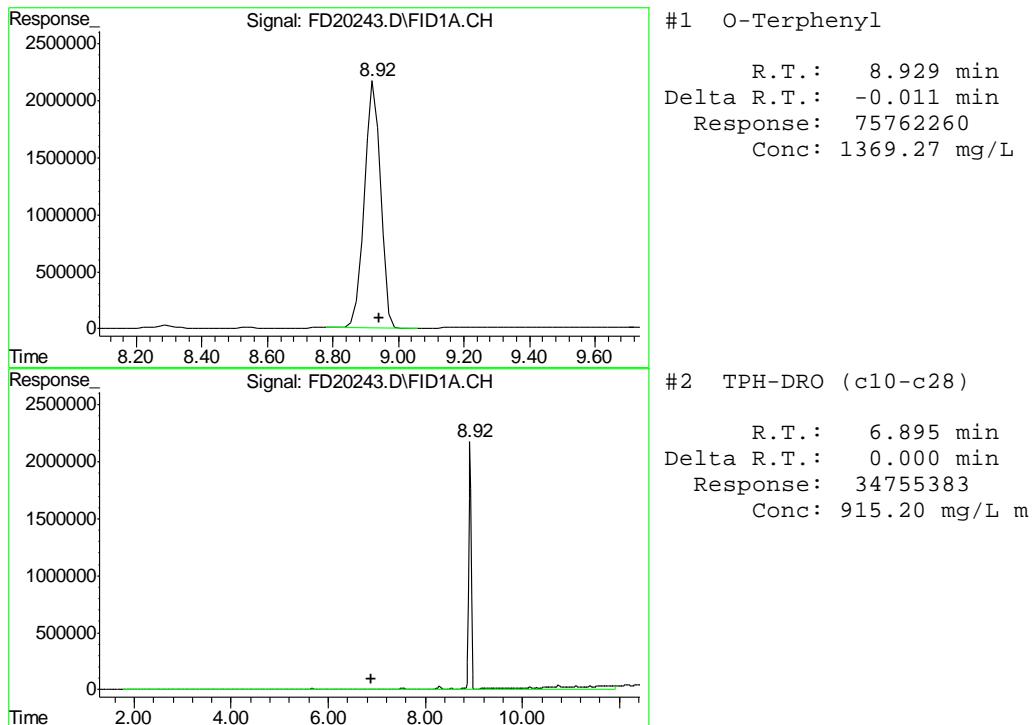
## Quantitation Report (QT Reviewed)

Data File : C:\MSDCHEM\2\DATA\2012\DEC\FD120712\FD20243.D Vial: 7  
 Acq On : 07 Dec 2012 11:12 am Operator: ashleyv  
 Sample : D41506-1 Inst : FID5  
 Misc : OP7061,GFD1015,30.00,,,1,1 Multiplr: 1.00  
 IntFile : autoint1.e  
 Quant Time: Dec 7 11:37 2012 Quant Results File: DRO-GFD982F.RES

Quant Method : C:\MSDCHEM\2...\DRO-GFD982F.M (Chemstation Integrator)  
 Title : 8015B TEH  
 Last Update : Fri Nov 16 10:24:56 2012  
 Response via : Multiple Level Calibration  
 DataAcq Meth : DRODUAL.M

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





## Quantitation Report (QT Reviewed)

Data File : C:\MSDCHEM\2\DATA\2012\DEC\FD120712\FD20235.D Vial: 3  
 Acq On : 12-7-2012 09:24:29 AM Operator: ashleyv  
 Sample : OP7061-MB Inst : FID5  
 Misc : OP7061,GFD1015,30.00,,,1,1 Multiplr: 1.00  
 IntFile : autoint1.e  
 Quant Time: Dec 07 10:09:28 2012 Quant Results File: DRO-GFD982F.RES

Quant Method : C:\MSDCHEM\2...\DRO-GFD982F.M (Chemstation Integrator)  
 Title : 8015B TEH  
 Last Update : Fri Nov 16 10:24:56 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
<hr/>				
System Monitoring Compounds				
1) S O-Terphenyl	8.93	86712214	1567.172	mg/L
<hr/>				
Target Compounds				
2) H TPH-DRO (c10-c28)	6.89	1719329	45.274	mg/L

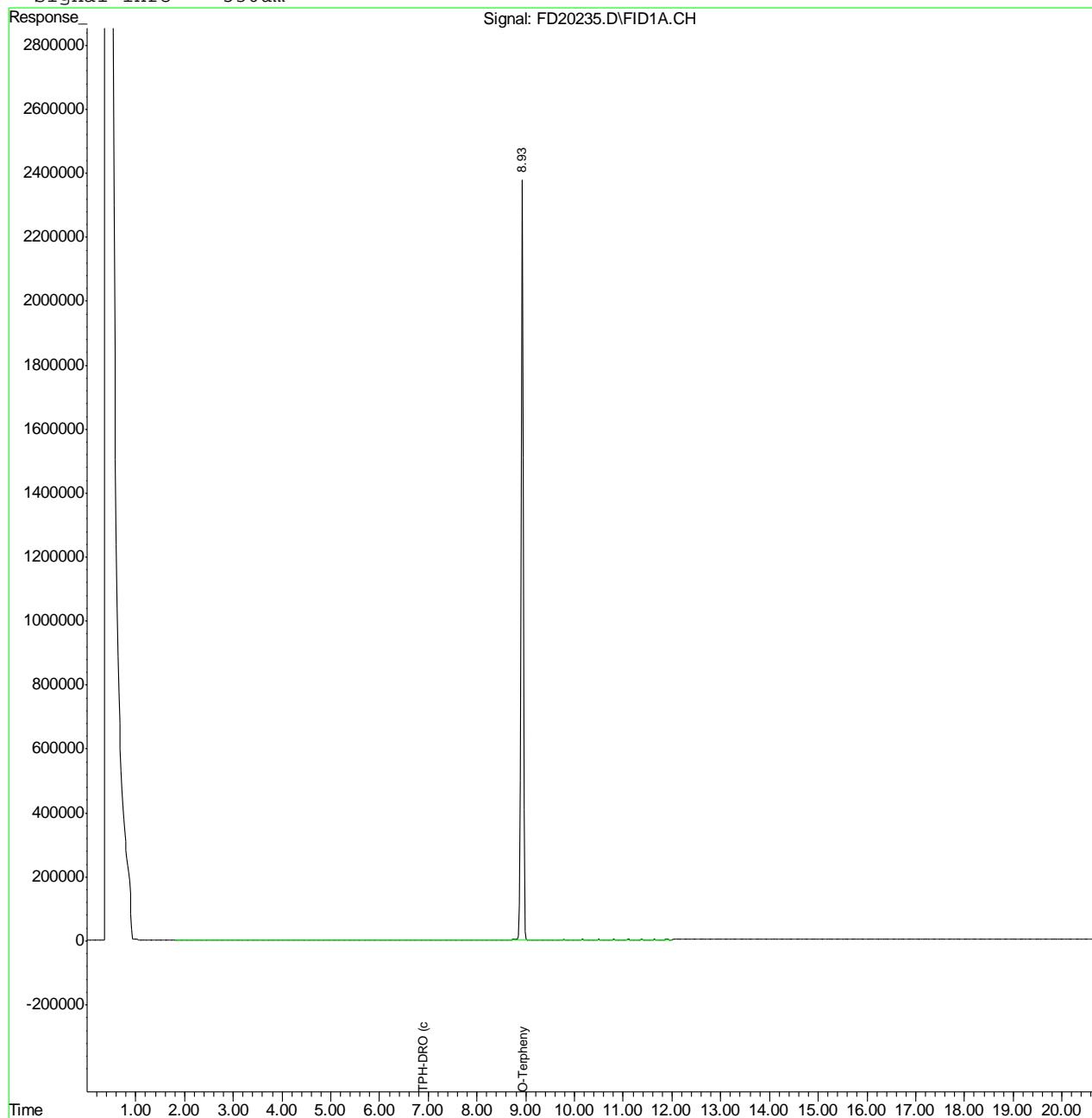
(f)=RT Delta > 1/2 Window (m)=manual int.  
 FD20235.D DRO-GFD982F.M Fri Dec 07 12:23:46 2012 GC

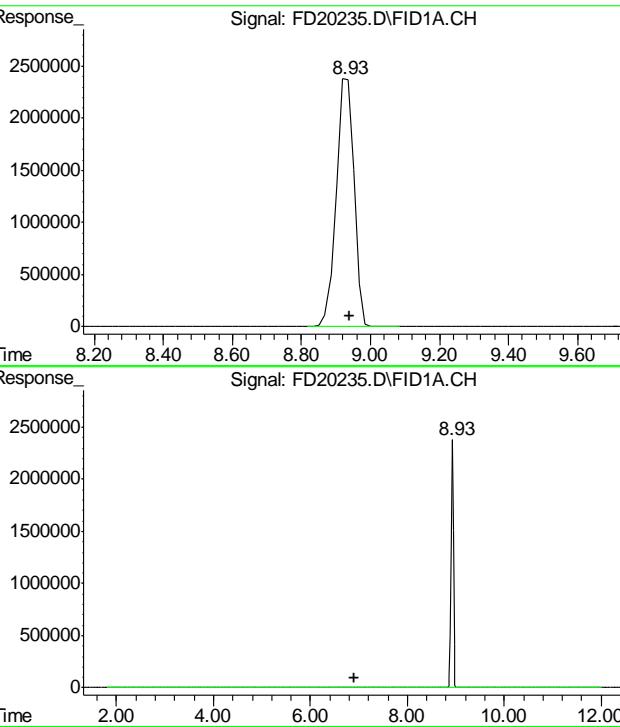
## Quantitation Report (QT Reviewed)

Data File : C:\MSDCHEM\2\DATA\2012\DEC\FD120712\FD20235.D Vial: 3  
 Acq On : 12-7-2012 09:24:29 AM Operator: ashleyv  
 Sample : OP7061-MB Inst : FID5  
 Misc : OP7061,GFD1015,30.00,,,1,1 Multiplr: 1.00  
 IntFile : autoint1.e  
 Quant Time: Dec 7 10:09 2012 Quant Results File: DRO-GFD982F.RES

Quant Method : C:\MSDCHEM\2...\DRO-GFD982F.M (Chemstation Integrator)  
 Title : 8015B TEH  
 Last Update : Fri Nov 16 10:24:56 2012  
 Response via : Multiple Level Calibration  
 DataAcq Meth : DRODUAL.M

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





#1 O-Terphenyl

R.T.: 8.934 min  
Delta R.T.: -0.006 min  
Response: 86712214  
Conc: 1567.17 mg/L

#2 TPH-DRO (c10-c28)

R.T.: 6.895 min  
Delta R.T.: 0.000 min  
Response: 1719329  
Conc: 45.27 mg/L



## 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: D41506  
Account: XTOKRWR - XTO Energy  
Project: PCU 296-5A

QC Batch ID: MP9008  
Matrix Type: SOLID

Methods: SW846 7471B  
Units: mg/kg

Prep Date:

12/06/12

Metal	RL	IDL	MDL	MB raw	final
Mercury	0.10	.0011	.0009	-0.00071	<0.10

Associated samples MP9008: D41506-1

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: D41506  
Account: XTOKRWR - XTO Energy  
Project: PCU 296-5A

QC Batch ID: MP9008  
Matrix Type: SOLID

Methods: SW846 7471B  
Units: mg/kg

Prep Date: 12/06/12

Metal	D41440-3 Original MS	Spikelot HGWSR1	QC % Rec	QC Limits
Mercury	0.26	0.96	0.806	86.9 75-125

Associated samples MP9008: D41506-1

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: D41506  
Account: XTOKRWR - XTO Energy  
Project: PCU 296-5A

QC Batch ID: MP9008  
Matrix Type: SOLID

Methods: SW846 7471B  
Units: mg/kg

Prep Date:

12/06/12

Metal	D41440-3 Original	MSD	Spikelot HGWSR1	MSD RPD	QC Limit
Mercury	0.26	1.0	0.819	90.4	4.1 20

Associated samples MP9008: D41506-1

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: D41506  
Account: XTOKRWR - XTO Energy  
Project: PCU 296-5A

QC Batch ID: MP9008  
Matrix Type: SOLID

Methods: SW846 7471B  
Units: mg/kg

Prep Date: 12/06/12

Metal	BSP Result	Spikelot HGWSR1	QC % Rec	QC Limits
Mercury	0.21	0.2	105.0	80-120

Associated samples MP9008: D41506-1

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

BLANK RESULTS SUMMARY  
Part 2 - Method Blanks

Login Number: D41506  
Account: XTOKRWR - XTO Energy  
Project: PCU 296-5A

QC Batch ID: MP9011  
Matrix Type: SOLID

Methods: SW846 6010C  
Units: mg/kg

Prep Date:

12/06/12

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	0.030	<1.0
Calcium	40	.54	9		
Chromium	1.0	.03	.03	0.050	<1.0
Cobalt	0.50	.04	.07		
Copper	1.0	.12	.15	0.080	<1.0
Iron	7.0	.12	.87		
Lead	5.0	.19	.24	0.38	<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	0.0	<3.0
Phosphorus	10	1.4	1.9		
Potassium	200	6.1	7		
Selenium	5.0	.48	.36	-0.070	<5.0
Silicon	5.0	.29	.37		
Silver	3.0	.04	.06	0.41	<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	1.1	<3.0

Associated samples MP9011: D41506-1

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

BLANK RESULTS SUMMARY  
Part 2 - Method Blanks

Login Number: D41506  
Account: XTOKRWR - XTO Energy  
Project: PCU 296-5A

QC Batch ID: MP9011  
Matrix Type: SOLID

Methods: SW846 6010C  
Units: mg/kg

Prep Date:

Metal

(anr) Analyte not requested

## MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D41506  
 Account: XTOKRWR - XTO Energy  
 Project: PCU 296-5A

QC Batch ID: MP9011  
 Matrix Type: SOLID

Methods: SW846 6010C  
 Units: mg/kg

Prep Date:

12/06/12

Metal	D41506-1 Original MS	Spikelot ICPALL2	% Rec	QC Limits
Aluminum	anr			
Antimony	anr			
Arsenic	anr			
Barium	2270	2320	222	22.6 (a) 75-125
Beryllium	anr			
Boron				
Cadmium	0.077	45.3	55.4	81.6 75-125
Calcium				
Chromium	26.9	68.1	55.4	80.6 75-125
Cobalt				
Copper	9.5	57.5	55.4	86.6 75-125
Iron	anr			
Lead	22.9	96.3	111	68.9N(b) 75-125
Lithium				
Magnesium	anr			
Manganese	anr			
Molybdenum	anr			
Nickel	14.1	57.7	55.4	82.6 75-125
Phosphorus	anr			
Potassium				
Selenium	0.56	95.3	111	85.5 75-125
Silicon				
Silver	0.0	20.1	22.2	90.7 75-125
Sodium				
Strontium				
Thallium	anr			
Tin				
Titanium				
Uranium	anr			
Vanadium				
Zinc	27.7	71.8	55.4	79.6 75-125

Associated samples MP9011: D41506-1

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

MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D41506  
Account: XTOKRWR - XTO Energy  
Project: PCU 296-5A

QC Batch ID: MP9011  
Matrix Type: SOLID

Methods: SW846 6010C  
Units: mg/kg

Prep Date:

Metal

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

14.2.2  
**14**

## MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D41506  
 Account: XTOKRWR - XTO Energy  
 Project: PCU 296-5A

QC Batch ID: MP9011  
 Matrix Type: SOLID

Methods: SW846 6010C  
 Units: mg/kg

Prep Date:

12/06/12

Metal	D41506-1 Original	MSD	Spikelot ICPALL2	% Rec	MSD RPD	QC Limit
Aluminum	anr					
Antimony	anr					
Arsenic	anr					
Barium	2270	4080	220	824.5(a)	55.0 (b)	20
Beryllium	anr					
Boron						
Cadmium	0.077	45.2	54.9	82.2	0.2	20
Calcium						
Chromium	26.9	67.7	54.9	80.7	0.6	20
Cobalt						
Copper	9.5	57.1	54.9	86.7	0.7	20
Iron	anr					
Lead	22.9	93.7	110	67.2N(c)	2.7	20
Lithium						
Magnesium	anr					
Manganese	anr					
Molybdenum	anr					
Nickel	14.1	52.3	54.9	73.6N(c)	9.8	20
Phosphorus	anr					
Potassium						
Selenium	0.56	95.7	110	86.7	0.4	20
Silicon						
Silver	0.0	20.0	22	91.1	0.5	20
Sodium						
Strontium						
Thallium	anr					
Tin						
Titanium						
Uranium	anr					
Vanadium						
Zinc	27.7	68.1	54.9	73.6N(c)	5.3	20

Associated samples MP9011: D41506-1

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

MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D41506  
Account: XTOKRWR - XTO Energy  
Project: PCU 296-5A

QC Batch ID: MP9011  
Matrix Type: SOLID

Methods: SW846 6010C  
Units: mg/kg

Prep Date:

Metal

- (N) Matrix Spike Rec. outside of QC limits
- (anr) Analyte not requested
- (a) Spike amount low relative to the sample amount. Refer to lab control or spike blank for recovery information.
- (b) High RPD due to possible sample matrix or nonhomogeneity.
- (c) Spike recovery indicates possible matrix interference.

14.2.2  
**14**

## SPIKE BLANK AND LAB CONTROL SAMPLE SUMMARY

Login Number: D41506  
 Account: XTOKRWR - XTO Energy  
 Project: PCU 296-5A

QC Batch ID: MP9011  
 Matrix Type: SOLID

Methods: SW846 6010C  
 Units: mg/kg

Prep Date: 12/06/12

Metal	BSP Result	Spikelot ICPALL2	% Rec	QC Limits
Aluminum	anr			
Antimony	anr			
Arsenic	anr			
Barium	213	200	106.5	80-120
Beryllium	anr			
Boron				
Cadmium	47.1	50	94.2	80-120
Calcium				
Chromium	50.6	50	101.2	80-120
Cobalt				
Copper	46.2	50	92.4	80-120
Iron	anr			
Lead	95.8	100	95.8	80-120
Lithium				
Magnesium	anr			
Manganese	anr			
Molybdenum	anr			
Nickel	47.7	50	95.4	80-120
Phosphorus	anr			
Potassium				
Selenium	97.3	100	97.3	80-120
Silicon				
Silver	20.2	20	101.0	80-120
Sodium				
Strontium				
Thallium	anr			
Tin				
Titanium				
Uranium	anr			
Vanadium				
Zinc	49.5	50	99.0	80-120

Associated samples MP9011: D41506-1

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

SPIKE BLANK AND LAB CONTROL SAMPLE SUMMARY

Login Number: D41506  
Account: XTOKRWR - XTO Energy  
Project: PCU 296-5A

QC Batch ID: MP9011  
Matrix Type: SOLID

Methods: SW846 6010C  
Units: mg/kg

Prep Date:

Metal

(anr) Analyte not requested

14.2.3  
**14**

## SERIAL DILUTION RESULTS SUMMARY

Login Number: D41506  
 Account: XTOKRWR - XTO Energy  
 Project: PCU 296-5A

QC Batch ID: MP9011  
 Matrix Type: SOLID

Methods: SW846 6010C  
 Units: ug/l

Prep Date: 12/06/12

Metal	D41506-1 Original	SDL 1:5	%DIF	QC Limits
Aluminum	anr			
Antimony	anr			
Arsenic	anr			
Barium	17300	20800	0.5	0-10
Beryllium	anr			
Boron				
Cadmium	0.00	0.00	NC (a)	0-10
Calcium				
Chromium	214	254	19.0*(b)	0-10
Cobalt				
Copper	86.5	91.0	5.2	0-10
Iron	anr			
Lead	209	217	19.6*(b)	0-10
Lithium				
Magnesium	anr			
Manganese	anr			
Molybdenum	anr			
Nickel	108	133	22.6*(b)	0-10
Phosphorus	anr			
Potassium				
Selenium	0.00	0.00	NC (a)	0-10
Silicon				
Silver	2.00	0.00	NC	0-10
Sodium				
Strontium				
Thallium	anr			
Tin				
Titanium				
Uranium	anr			
Vanadium				
Zinc	252	326	29.2*(b)	0-10

Associated samples MP9011: D41506-1

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

SERIAL DILUTION RESULTS SUMMARY

Login Number: D41506  
Account: XTOKRWR - XTO Energy  
Project: PCU 296-5A

QC Batch ID: MP9011  
Matrix Type: SOLID

Methods: SW846 6010C  
Units: ug/l

Prep Date:

Metal

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

14.2.4  
**14**

BLANK RESULTS SUMMARY  
Part 2 - Method Blanks

Login Number: D41506  
Account: XTOKRWR - XTO Energy  
Project: PCU 296-5A

QC Batch ID: MP9012  
Matrix Type: SOLID

Methods: SW846 6020A  
Units: mg/kg

Prep Date:

12/06/12

Metal	RL	IDL	MDL	MB raw	final
Aluminum	25	.22	.31		
Antimony	0.20	.0018	.0075		
Arsenic	0.10	.006	.06	-0.0052	<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 MP9012: D41506-1

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

14.3.1  
14

## MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D41506  
 Account: XTOKWR - XTO Energy  
 Project: PCU 296-5A

QC Batch ID: MP9012  
 Matrix Type: SOLID

Methods: SW846 6020A  
 Units: mg/kg

Prep Date:

12/06/12

Metal	D41506-1 Original MS	Spikelot ICPALL2	% Rec	QC Limits
Aluminum				
Antimony				
Arsenic	5.0	97.5	111	83.4    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 MP9012: D41506-1

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: D41506  
 Account: XTOKRWR - XTO Energy  
 Project: PCU 296-5A

QC Batch ID: MP9012  
 Matrix Type: SOLID

Methods: SW846 6020A  
 Units: mg/kg

Prep Date:

12/06/12

Metal	D41506-1 Original	MSD	Spikelot ICPALL2	% Rec	MSD RPD	QC Limit
Aluminum						
Antimony						
Arsenic	5.0	97.5	110	84.3	0.0	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 MP9012: D41506-1

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: D41506  
 Account: XTOKRWR - XTO Energy  
 Project: PCU 296-5A

QC Batch ID: MP9012  
 Matrix Type: SOLID

Methods: SW846 6020A  
 Units: mg/kg

Prep Date: 12/06/12

Metal	BSP Result	Spikelot ICPALL2	% Rec	QC Limits
Aluminum				
Antimony				
Arsenic	84.7	100	84.7	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 MP9012: D41506-1

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

14.3.3  
**14**

## SERIAL DILUTION RESULTS SUMMARY

Login Number: D41506  
 Account: XTOKRWR - XTO Energy  
 Project: PCU 296-5A

QC Batch ID: MP9012  
 Matrix Type: SOLID

Methods: SW846 6020A  
 Units: ug/l

Prep Date:

12/06/12

Metal	D41506-1	Original	SDL	5:25 %DIF	QC Limits
-------	----------	----------	-----	-----------	--------------

Aluminum  
 Antimony  
 Arsenic 45.4 42.8 5.6 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 MP9012: D41506-1

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

14.3.4  
14

BLANK RESULTS SUMMARY  
Part 2 - Method Blanks

Login Number: D41506  
Account: XTOKRWR - XTO Energy  
Project: PCU 296-5A

QC Batch ID: MP9015  
Matrix Type: AQUEOUS

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

Prep Date:

12/06/12

Metal	RL	IDL	MDL	MB raw	final
Aluminum	500	48	130		
Antimony	150	8.5	18		
Arsenic	130	22	42		
Barium	50	.5	9		
Beryllium	50	6.5	16		
Boron	250	5	22		
Cadmium	50	3	3		
Calcium	2000	27	80	9.5	<2000
Chromium	50	1.5	2.8		
Cobalt	25	2	2.1		
Copper	50	6	15		
Iron	350	6	100		
Lead	250	9.5	15		
Lithium	10	2.5			
Magnesium	1000	33	110	26.0	<1000
Manganese	25	6	6		
Molybdenum	50	11	11		
Nickel	150	2.5	2.9		
Phosphorus	500	70	300		
Potassium	5000	310	750		
Selenium	250	24	55		
Silicon	250	15			
Silver	150	2	4.9		
Sodium	2000	30	490	103	<2000
Strontium	25	.2	7.5		
Thallium	50	15	43		
Tin	250	60			
Titanium	50	.5			
Uranium	250	11	23		
Vanadium	50	1	2.4		
Zinc	150	2.5	12		

Associated samples MP9015: D41506-1A

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

BLANK RESULTS SUMMARY  
Part 2 - Method Blanks

Login Number: D41506  
Account: XTOKRWR - XTO Energy  
Project: PCU 296-5A

QC Batch ID: MP9015  
Matrix Type: AQUEOUS

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

Prep Date:

Metal

(anr) Analyte not requested

## MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D41506  
 Account: XTOKRWR - XTO Energy  
 Project: PCU 296-5A

QC Batch ID: MP9015  
 Matrix Type: AQUEOUS

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

Prep Date: 12/06/12

Metal	D41381-1A Original MS	Spikelot ICPALL2	% Rec	QC Limits
Aluminum				
Antimony				
Arsenic				
Barium				
Beryllium				
Boron				
Cadmium				
Calcium	32600	168000	125000	108.3
Chromium				
Cobalt				
Copper				
Iron				
Lead				
Lithium				
Magnesium	8180	132000	125000	99.1
Manganese				
Molybdenum				
Nickel				
Phosphorus				
Potassium				
Selenium				
Silicon				
Silver				
Sodium	35000	161000	125000	100.8
Strontium				
Thallium				
Tin				
Titanium				
Uranium				
Vanadium				
Zinc				

Associated samples MP9015: D41506-1A

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

MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D41506  
Account: XTOKRWR - XTO Energy  
Project: PCU 296-5A

QC Batch ID: MP9015  
Matrix Type: AQUEOUS

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

Prep Date:

Metal

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

## MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D41506  
 Account: XTOKRWR - XTO Energy  
 Project: PCU 296-5A

QC Batch ID: MP9015  
 Matrix Type: AQUEOUS

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

Prep Date: 12/06/12

Metal	D41381-1A Original MSD	Spikelot ICPALL2	MSD % Rec	MSD RPD	QC Limit
Aluminum					
Antimony					
Arsenic					
Barium					
Beryllium					
Boron					
Cadmium					
Calcium	32600	169000	125000	109.1	0.6
Chromium					
Cobalt					
Copper					
Iron					
Lead					
Lithium					
Magnesium	8180	132000	125000	99.1	0.0
Manganese					
Molybdenum					
Nickel					
Phosphorus					
Potassium					
Selenium					
Silicon					
Silver					
Sodium	35000	161000	125000	100.8	0.0
Strontium					
Thallium					
Tin					
Titanium					
Uranium					
Vanadium					
Zinc					

Associated samples MP9015: D41506-1A

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

MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D41506  
Account: XTOKRWR - XTO Energy  
Project: PCU 296-5A

QC Batch ID: MP9015  
Matrix Type: AQUEOUS

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

Prep Date:

Metal

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

## SPIKE BLANK AND LAB CONTROL SAMPLE SUMMARY

Login Number: D41506  
 Account: XTOKRWR - XTO Energy  
 Project: PCU 296-5A

QC Batch ID: MP9015  
 Matrix Type: AQUEOUS

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

Prep Date: 12/06/12

Metal	BSP Result	Spikelot ICPALL2	% Rec	QC Limits
Aluminum				
Antimony				
Arsenic				
Barium				
Beryllium				
Boron				
Cadmium				
Calcium	135000	125000	108.0	80-120
Chromium				
Cobalt				
Copper				
Iron				
Lead				
Lithium				
Magnesium	122000	125000	97.6	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 MP9015: D41506-1A

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

SPIKE BLANK AND LAB CONTROL SAMPLE SUMMARY

Login Number: D41506  
Account: XTOKRWR - XTO Energy  
Project: PCU 296-5A

QC Batch ID: MP9015  
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: D41506  
 Account: XTOKRWR - XTO Energy  
 Project: PCU 296-5A

QC Batch ID: MP9015  
 Matrix Type: AQUEOUS

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

Prep Date: 12/06/12

Metal	D41381-1A	Original	SDL 1:5	%DIF	QC Limits
Aluminum					
Antimony					
Arsenic					
Barium					
Beryllium					
Boron					
Cadmium					
Calcium	6510	6470	0.8		0-10
Chromium					
Cobalt					
Copper					
Iron					
Lead					
Lithium					
Magnesium	1640	1720	4.8		0-10
Manganese					
Molybdenum					
Nickel					
Phosphorus					
Potassium					
Selenium					
Silicon					
Silver					
Sodium	7000	7390	5.5		0-10
Strontium					
Thallium					
Tin					
Titanium					
Uranium					
Vanadium					
Zinc					

Associated samples MP9015: D41506-1A

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

SERIAL DILUTION RESULTS SUMMARY

Login Number: D41506  
Account: XTOKRWR - XTO Energy  
Project: PCU 296-5A

QC Batch ID: MP9015  
Matrix Type: AQUEOUS

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

Prep Date:

Metal

(anr) Analyte not requested



## General Chemistry

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### QC Data Summaries

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Includes the following where applicable:

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

METHOD BLANK AND SPIKE RESULTS SUMMARY  
GENERAL CHEMISTRY

Login Number: D41506  
Account: XTOKWR - XTO Energy  
Project: PCU 296-5A

Analyte	Batch ID	RL	MB Result	Units	Spike Amount	BSP Result	BSP %Recov	QC Limits
Chromium, Hexavalent	GP8845/GN17994	1.0	0.0	mg/kg	176.0	170	97.0	80-120%
Specific Conductivity	GP8856/GN18003			umhos/cm	9992	10000	100.5	90-110%
pH	GN17963			su	8.00	7.99	99.9	99.3-100.7%

Associated Samples:

Batch GP8845: D41506-1

Batch GP8856: D41506-1

Batch GN17963: D41506-1

(\*) Outside of QC limits

DUPLICATE RESULTS SUMMARY  
GENERAL CHEMISTRY

Login Number: D41506  
Account: XTOKWR - XTO Energy  
Project: PCU 296-5A

Analyte	Batch ID	QC Sample	Units	Original Result	DUP Result	RPD	QC Limits
Chromium, Hexavalent	GP8845/GN17994	D41506-1	mg/kg	0.0	0.0	0.0	0-20%

Associated Samples:  
Batch GP8845: D41506-1  
(\*) Outside of QC limits

MATRIX SPIKE RESULTS SUMMARY  
GENERAL CHEMISTRY

Login Number: D41506  
Account: XTOKWR - XTO Energy  
Project: PCU 296-5A

Analyte	Batch ID	QC Sample	Units	Original Result	Spike Amount	MS Result	%Rec	QC Limits
Chromium, Hexavalent	GP8845/GN17994	D41506-1	mg/kg	0.0	40.0	35.7	89.2	75-125%

Associated Samples:

Batch GP8845: D41506-1

(\*) Outside of QC limits

(N) Matrix Spike Rec. outside of QC limits

MATRIX SPIKE DUPLICATE RESULTS SUMMARY  
GENERAL CHEMISTRY

Login Number: D41506  
Account: XTOKWR - XTO Energy  
Project: PCU 296-5A

Analyte	Batch ID	QC Sample	Units	Original Result	Spike Amount	MSD Result	RPD	QC Limit
Chromium, Hexavalent	GP8845/GN17994	D41506-1	mg/kg	0.0	40.0	38.7	7.9	20%

Associated Samples:

Batch GP8845: D41506-1

(\*) Outside of QC limits

(N) Matrix Spike Rec. outside of QC limits