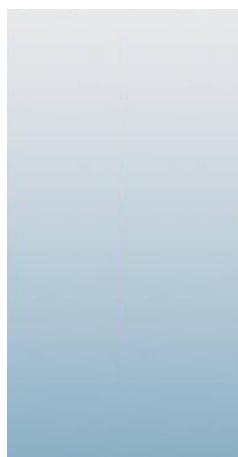




12/12/12



Technical Report for

XTO Energy

PCU 296-5A

1210-04

Accutest Job Number: D41448

Sampling Date: 11/30/12

Report to:

KRW Consulting, Inc.
8000 West 14th Avenue
Lakewood, CO 80214
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ATTN: Dwayne Knudson

Total number of pages in report: 146



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: D41448

PCU 296-5A

Project No: 1210-04

Sample Number	Collected Date	Time By	Received	Matrix Code	Type	Client Sample ID
D41448-1	11/30/12	11:15 DK	12/04/12	SO	Soil	RP POST SOLIDIFICATION
D41448-1A	11/30/12	11:15 DK	12/04/12	SO	Soil	RP POST SOLIDIFICATION

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



CASE NARRATIVE / CONFORMANCE SUMMARY

Client: XTO Energy

Job No D41448

Site: PCU 296-5A

Report Date 12/10/2012 4:14:27 PM

On 12/04/2012, 1 sample(s), 0 Trip Blank(s), and 0 Field Blank(s) were received at Accutest Mountain States (AMS) at a temperature of 2.8 °C. The samples were intact and properly preserved, unless noted below. An AMS Job Number of D41448 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

Matrix SO	Batch ID: V3V1284
------------------	--------------------------

- All samples were analyzed within the recommended method holding time.
- Sample(s) D41448-1MS, D41448-1MSD 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

Matrix SO	Batch ID: 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

Matrix SO	Batch ID: GGB1023
------------------	--------------------------

- 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 were used as the QC samples indicated.

Extractables by GC By Method SW846-8015B

Matrix SO	Batch ID: OP7053
------------------	-------------------------

- 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.

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: MP9006

- 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) D41381-1MS, D41381-1MSD, D41381-1SDL were used as the QC samples for the metals analysis.
- The matrix spike (MS) recovery(s) of Barium, Cadmium, Lead, Nickel are outside control limits. Spike recovery indicates possible matrix interference and/or sample nonhomogeneity.
- The matrix spike duplicate (MSD) recovery(s) of Cadmium, Lead, Nickel are outside control limits. Probable cause due to matrix interference.
- The serial dilution RPD(s) for Cadmium, Silver, Chromium, Nickel, Zinc are outside control limits for sample MP9006-SD1. Percent difference acceptable due to low initial sample concentration (< 50 times IDL).
- MP9006-MB1 for Barium: All sample results >10x method blank concentration or <RL.
- MP9006-S1 for Nickel: Spike recovery indicates possible matrix interference.
- MP9006-SD1 for Zinc: Serial dilution indicates possible matrix interference.
- MP9006-SD1 for Nickel: Serial dilution indicates possible matrix interference.
- MP9006-SD1 for Chromium: Serial dilution indicates possible matrix interference.
- MP9006-S1 for Cadmium: Spike recovery indicates possible matrix interference.
- MP9006-S1 for Lead: Spike recovery indicates possible matrix interference.

Metals By Method SW846 6020A

Matrix SO

Batch ID: MP9007

- 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) D41381-1MS, D41381-1MSD, D41381-1SDL 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: GN17944

- 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.
- D41448-1 for Chromium, Hexavalent: Dilution required due to matrix interference.

Wet Chemistry By Method SW846 3060A/7196A M

Matrix SO

Batch ID: R15368

- The data for SW846 3060A/7196A M meets quality control requirements.
- D41448-1 for Chromium, Trivalent: Calculated as: (Chromium) - (Chromium, Hexavalent)

Wet Chemistry By Method SW846 9045D

Matrix SO

Batch ID: GN17962

- The following samples were run outside of holding time for method SW846 9045D: D41448-1

Wet Chemistry By Method USDA HANDBOOK 60

Matrix SO

Batch ID: MP9015

- D41448-1A 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: D41448
Account: XTO Energy
Project: PCU 296-5A
Collected: 11/30/12

3

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

D41448-1 RP POST SOLIDIFICATION

Benzene	0.128	0.11	0.056	mg/kg	SW846 8260B
Toluene	0.327	0.22	0.11	mg/kg	SW846 8260B
Ethylbenzene	0.0466 J	0.22	0.042	mg/kg	SW846 8260B
Xylene (total)	0.399 J	0.45	0.22	mg/kg	SW846 8260B
Chrysene	0.0107 J	0.013	0.0070	mg/kg	SW846 8270C BY SIM
Fluorene	0.0149	0.013	0.0070	mg/kg	SW846 8270C BY SIM
Naphthalene	0.0576	0.019	0.017	mg/kg	SW846 8270C BY SIM
Pyrene	0.0080 J	0.013	0.0070	mg/kg	SW846 8270C BY SIM
TPH-GRO (C6-C10)	12.3 J	22	11	mg/kg	SW846 8015B
TPH-DRO (C10-C28)	159	11	6.5	mg/kg	SW846-8015B
Arsenic	9.9	0.16		mg/kg	SW846 6020A
Barium	11400	16		mg/kg	SW846 6010C
Chromium	14.1	1.6		mg/kg	SW846 6010C
Copper	28.4	1.6		mg/kg	SW846 6010C
Lead	15.5	8.1		mg/kg	SW846 6010C
Nickel	112	4.9		mg/kg	SW846 6010C
Zinc	33.8	4.9		mg/kg	SW846 6010C
Specific Conductivity	8770	1.0		umhos/cm	SM2510B-1997 MOD
Chromium, Trivalent ^a	14.1	6.6		mg/kg	SW846 3060A/7196A M
Redox Potential Vs H2	173			mv	ASTM D1498-76M
pH	12.32			su	SW846 9045D

D41448-1A RP POST SOLIDIFICATION

Calcium	393	2.0	mg/l	SW846 6010C
Sodium	981	2.0	mg/l	SW846 6010C
Sodium Adsorption Ratio ^b	13.6		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

Report of Analysis

Accutest Laboratories

Report of Analysis

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Client Sample ID: RP POST SOLIDIFICATION**Lab Sample ID:** D41448-1**Matrix:** SO - Soil**Method:** SW846 8260B**Project:** PCU 296-5A**Date Sampled:** 11/30/12**Date Received:** 12/04/12**Percent Solids:** 61.7

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	3V21984.D	1	12/06/12	BD	n/a	n/a	V3V1284
Run #2							

	Initial Weight	Final Volume	Methanol Aliquot
Run #1	5.02 g	5.0 ml	100 ul
Run #2			

Purgeable Aromatics

CAS No.	Compound	Result	RL	MDL	Units	Q
71-43-2	Benzene	0.128	0.11	0.056	mg/kg	
108-88-3	Toluene	0.327	0.22	0.11	mg/kg	
100-41-4	Ethylbenzene	0.0466	0.22	0.042	mg/kg	J
1330-20-7	Xylene (total)	0.399	0.45	0.22	mg/kg	J

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
2037-26-5	Toluene-D8	98%		64-130%
460-00-4	4-Bromofluorobenzene	94%		62-131%
17060-07-0	1,2-Dichloroethane-D4	111%		70-130%

ND = Not detected MDL - Method Detection Limit

J = Indicates an estimated value

RL = Reporting Limit

B = Indicates analyte found in associated method blank

E = Indicates value exceeds calibration range

N = Indicates presumptive evidence of a compound

Accutest Laboratories

Report of Analysis

Page 1 of 1

Client Sample ID:	RP POST SOLIDIFICATION	Date Sampled:	11/30/12
Lab Sample ID:	D41448-1	Date Received:	12/04/12
Matrix:	SO - Soil	Percent Solids:	61.7
Method:	SW846 8270C BY SIM	SW846 3546	
Project:	PCU 296-5A		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	3G12515.D	1	12/10/12	DC	12/10/12	OP7075	E3G593
Run #2							

	Initial Weight	Final Volume
Run #1	30.0 g	1.0 ml
Run #2		

COGCC Table 910-1 PAH List

CAS No.	Compound	Result	RL	MDL	Units	Q
83-32-9	Acenaphthene	ND	0.013	0.0070	mg/kg	
120-12-7	Anthracene	ND	0.013	0.0070	mg/kg	
56-55-3	Benzo(a)anthracene	ND	0.013	0.0070	mg/kg	
205-99-2	Benzo(b)fluoranthene	ND	0.013	0.0070	mg/kg	
207-08-9	Benzo(k)fluoranthene	ND	0.013	0.0070	mg/kg	
50-32-8	Benzo(a)pyrene	ND	0.013	0.0070	mg/kg	
218-01-9	Chrysene	0.0107	0.013	0.0070	mg/kg	J
53-70-3	Dibenz(a,h)anthracene	ND	0.013	0.0070	mg/kg	
206-44-0	Fluoranthene	ND	0.013	0.0070	mg/kg	
86-73-7	Fluorene	0.0149	0.013	0.0070	mg/kg	
193-39-5	Indeno(1,2,3-cd)pyrene	ND	0.013	0.0070	mg/kg	
91-20-3	Naphthalene	0.0576	0.019	0.017	mg/kg	
129-00-0	Pyrene	0.0080	0.013	0.0070	mg/kg	J

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
4165-60-0	Nitrobenzene-d5	53%		10-159%
321-60-8	2-Fluorobiphenyl	50%		19-131%
1718-51-0	Terphenyl-d14	69%		18-150%

ND = Not detected MDL - Method Detection Limit

J = Indicates an estimated value

RL = Reporting Limit

B = Indicates analyte found in associated method blank

E = Indicates value exceeds calibration range

N = Indicates presumptive evidence of a compound

4.1

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Accutest Laboratories

Report of Analysis

Page 1 of 1

Client Sample ID: RP POST SOLIDIFICATION**Lab Sample ID:** D41448-1**Date Sampled:** 11/30/12**Matrix:** SO - Soil**Date Received:** 12/04/12**Method:** SW846 8015B**Percent Solids:** 61.7**Project:** PCU 296-5A

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	GB18767.D	1	12/06/12	SK	n/a	n/a	GGB1023
Run #2							

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

CAS No.	Compound	Result	RL	MDL	Units	Q
	TPH-GRO (C6-C10)	12.3	22	11	mg/kg	J
CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits		
120-82-1	1,2,4-Trichlorobenzene	83%		60-140%		

ND = Not detected MDL - Method Detection Limit

J = Indicates an estimated value

RL = Reporting Limit

B = Indicates analyte found in associated method blank

E = Indicates value exceeds calibration range

N = Indicates presumptive evidence of a compound

4.1

4

Accutest Laboratories

Report of Analysis

Page 1 of 1

Client Sample ID: RP POST SOLIDIFICATION**Lab Sample ID:** D41448-1**Date Sampled:** 11/30/12**Matrix:** SO - Soil**Date Received:** 12/04/12**Method:** SW846-8015B SW846 3546**Percent Solids:** 61.7**Project:** PCU 296-5A

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	FH008082.D	1	12/06/12	TR	12/05/12	OP7053	GFH446
Run #2							

	Initial Weight	Final Volume
Run #1	30.1 g	1.0 ml
Run #2		

CAS No.	Compound	Result	RL	MDL	Units	Q
	TPH-DRO (C10-C28)	159	11	6.5	mg/kg	
CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits		
84-15-1	o-Terphenyl	64%		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 POST SOLIDIFICATION	Date Sampled:	11/30/12
Lab Sample ID:	D41448-1	Date Received:	12/04/12
Matrix:	SO - Soil	Percent Solids:	61.7
Project:	PCU 296-5A		

Metals Analysis

Analyte	Result	RL	Units	DF	Prep	Analyzed By	Method	Prep Method
Arsenic	9.9	0.16	mg/kg	5	12/05/12	12/07/12 JM	SW846 6020A ³	SW846 3050B ⁶
Barium	11400	16	mg/kg	10	12/05/12	12/07/12 JB	SW846 6010C ⁴	SW846 3050B ⁵
Cadmium	< 1.6	1.6	mg/kg	1	12/05/12	12/06/12 JB	SW846 6010C ¹	SW846 3050B ⁵
Chromium	14.1	1.6	mg/kg	1	12/05/12	12/06/12 JB	SW846 6010C ¹	SW846 3050B ⁵
Copper	28.4	1.6	mg/kg	1	12/05/12	12/06/12 JB	SW846 6010C ¹	SW846 3050B ⁵
Lead	15.5	8.1	mg/kg	1	12/05/12	12/06/12 JB	SW846 6010C ¹	SW846 3050B ⁵
Mercury	< 0.13	0.13	mg/kg	1	12/06/12	12/06/12 JB	SW846 7471B ²	SW846 7471B ⁷
Nickel	112	4.9	mg/kg	1	12/05/12	12/06/12 JB	SW846 6010C ¹	SW846 3050B ⁵
Selenium	< 8.1	8.1	mg/kg	1	12/05/12	12/06/12 JB	SW846 6010C ¹	SW846 3050B ⁵
Silver	< 4.9	4.9	mg/kg	1	12/05/12	12/06/12 JB	SW846 6010C ¹	SW846 3050B ⁵
Zinc	33.8	4.9	mg/kg	1	12/05/12	12/06/12 JB	SW846 6010C ¹	SW846 3050B ⁵

- (1) Instrument QC Batch: MA3055
- (2) Instrument QC Batch: MA3057
- (3) Instrument QC Batch: MA3061
- (4) Instrument QC Batch: MA3062
- (5) Prep QC Batch: MP9006
- (6) Prep QC Batch: MP9007
- (7) Prep QC Batch: MP9008

RL = Reporting Limit

Report of Analysis

Page 1 of 1

Client Sample ID: RP POST SOLIDIFICATION**Lab Sample ID:** D41448-1**Matrix:** SO - Soil**Project:** PCU 296-5A**Date Sampled:** 11/30/12**Date Received:** 12/04/12**Percent Solids:** 61.7**General Chemistry**

Analyte	Result	RL	Units	DF	Analyzed	By	Method
prep: DEPT.OF AG, BOOK N9							
Specific Conductivity	8770	1.0	umhos/cm	1	12/07/12	JD	SM2510B-1997 MOD
Chromium, Hexavalent ^a	< 5.0	5.0	mg/kg	5	12/07/12	KB	SW846 3060A/7196A
Chromium, Trivalent ^b	14.1	6.6	mg/kg	1	12/07/12	KB	SW846 3060A/7196A M
Redox Potential Vs H2	173		mv	1	12/05/12	JD	ASTM D1498-76M
Solids, Percent	61.7		%	1	12/05/12	SWT	SM19 2540B M
pH	12.32		su	1	12/05/12 15:00	JD	SW846 9045D

(a) Dilution required due to matrix interference.

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

RL = Reporting Limit

Report of Analysis

Page 1 of 1

Client Sample ID: RP POST SOLIDIFICATION**Lab Sample ID:** D41448-1A**Matrix:** SO - Soil**Project:** PCU 296-5A**Date Sampled:** 11/30/12**Date Received:** 12/04/12**Percent Solids:** 61.7**SAR Metals Analysis**

Analyte	Result	RL	Units	DF	Prep	Analyzed By	Method	Prep Method
Calcium	393	2.0	mg/l	1	12/06/12	12/10/12 JB	SW846 6010C ¹	SW846 3010A/M ²
Magnesium	< 1.0	1.0	mg/l	1	12/06/12	12/10/12 JB	SW846 6010C ¹	SW846 3010A/M ²
Sodium	981	2.0	mg/l	1	12/06/12	12/10/12 JB	SW846 6010C ¹	SW846 3010A/M ²

(1) Instrument QC Batch: MA3069

(2) Prep QC Batch: MP9015

RL = Reporting Limit

Report of Analysis

Page 1 of 1

Client Sample ID: RP POST SOLIDIFICATION**Lab Sample ID:** D41448-1A**Matrix:** SO - Soil**Project:** PCU 296-5A**Date Sampled:** 11/30/12**Date Received:** 12/04/12**Percent Solids:** 61.7**General Chemistry**

Analyte	Result	RL	Units	DF	Analyzed	By	Method
Sodium Adsorption Ratio ^a	13.6		ratio	1	12/10/12 10:06	JB	USDA HANDBOOK 60

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

RL = Reporting Limit



Misc. Forms

5

Custody Documents and Other Forms

Includes the following where applicable:

- Chain of Custody



CHAIN OF CUSTODY

PAGE 1 OF 1

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

FED-EX Tracking #	Bottle Order Control #
Accutest Job #	D41448

Client / Reporting Information		Project Information		Requested Analysis (see TEST CODE sheet)										Matrix Codes			
Company Name KRW Consulting	Project Name: XTO P2296 - SA																
Street Address 8000 West 14th Street; Suite 200	Street:			Billing Information (If different from Report to)													
City Lakewood, CO 80214	City	State	Company Name XTO Energy														
Project Contact Dwayne Knudson	Project # 1210-04			Street Address 21459 CR 5													
Phone # 970-488-1098	Client Purchase Order # Joe Hess			City Rifle, CO 81650													
Sampler(s) Name(s) DWAYNE KNUDSON	Project Manager Jessica Dooling			Attention: T - 910													
				Number of preserved Bottles													
Accutest Sample #	Field ID / Point of Collection	MECH/DI Vial #	Date	Time	Sampled by	Matrix	# of bottles	HCl	NaOH	HFCA	None	Di Water	MECH	ENCORE	Bottles		
	RP POST SOLIDIFICATION		11-30-12	11:15	DK	SO	5			X						T - 910	
Turnaround Time (Business days)		Approved By (Accutest PM): / Date:		Data Deliverable Information										Comments / Special Instructions			
<input type="checkbox"/> Std. 10 Business Days	<input type="checkbox"/> Std. 5 Business Days (By contract only)			<input type="checkbox"/> Commercial "A" (Level 1)	<input type="checkbox"/> State Forms Required												
<input type="checkbox"/> 3 Day Emergency	<input type="checkbox"/> 2 Day Emergency			<input type="checkbox"/> Commercial "B" (Level 2)	<input type="checkbox"/> Send Forms to State												
<input type="checkbox"/> 1 Day Emergency	<input type="checkbox"/>			<input type="checkbox"/> COMMNB	<input type="checkbox"/> Report by Fax												
Emergency & Rush T/A data available VIA Lablink														<input checked="" type="checkbox"/> COMMNB+	<input checked="" type="checkbox"/> Report by PDF ONLY		
														<input type="checkbox"/> EDD Format			
														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: 1 Tom Albranson	Date Time: 11/3/12 15:30	Received By: 1	Relinquished By: 2	Date Time: 12/1/12 12:00	Received By: 2												
Relinquished by Sampler: 3	Date Time: 11/3/12 15:30	Received By: 3	Relinquished By: 4	Date Time: 12/1/12 12:00	Received By: 4												
Relinquished by: 5	Date Time: 11/3/12 15:30	Received By: 5	Custody Seal # HDCC	<input checked="" type="checkbox"/> Intact	Preserved where applicable	On Ice	Cooler Temp.										

D41448: Chain of Custody

Page 1 of 2



Accutest Laboratories Sample Receipt Summary

Accutest Job Number: D41448

Client: KRW CONSULTING

Immediate Client Services Action Required: No

Date / Time Received: 12/4/2012 12:20:00 PM

No. Coolers:

1

Client Service Action Required at Login: No

Project: XTO PCU296-5A

Airbill #'s: HDCO

Cooler Security Y or N

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

Cooler Temperature Y or N

- | | | |
|------------------------------|-------------------------------------|--------------------------|
| 1. Temp criteria achieved: | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| 2. Cooler temp verification: | Infrared gun | |
| 3. Cooler media: | Ice (bag) | |

Quality Control Preservation Y or N N/A

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

Sample Integrity - Documentation

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

Sample Integrity - Condition

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

Sample Integrity - Instructions

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

Comments

Accutest Laboratories
V:(303) 425-6021

4036 Youngfield Street
F: (303) 425-6854

Wheat Ridge, CO
www.accutest.com

5.1

5

D41448: Chain of Custody

Page 2 of 2



GC/MS Volatiles

QC Data Summaries

Includes the following where applicable:

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



Method Blank Summary

Page 1 of 1

Job Number: D41448
Account: XTOKRWR XTO Energy
Project: PCU 296-5A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
V3V1284-MB	3V21977.D	1	12/06/12	BD	n/a	n/a	V3V1284

The QC reported here applies to the following samples:

Method: SW846 8260B

D41448-1

6.1

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

CAS No.	Surrogate Recoveries	Limits
2037-26-5	Toluene-D8	103%
460-00-4	4-Bromofluorobenzene	86%
17060-07-0	1,2-Dichloroethane-D4	118%

Blank Spike Summary

Job Number: D41448
Account: XTOKWR XTO Energy
Project: PCU 296-5A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
V3V1284-BS	3V21981.D	1	12/06/12	BD	n/a	n/a	V3V1284

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

D41448-1

CAS No.	Compound	Spike ug/kg	BSP ug/kg	BSP %	Limits
71-43-2	Benzene	50	48.3	97	70-130
100-41-4	Ethylbenzene	50	46.4	93	70-130
108-88-3	Toluene	50	44.5	89	70-130
1330-20-7	Xylene (total)	150	139	93	70-130

CAS No.	Surrogate Recoveries	BSP	Limits
2037-26-5	Toluene-D8	99%	64-130%
460-00-4	4-Bromofluorobenzene	94%	62-131%
17060-07-0	1,2-Dichloroethane-D4	114%	70-130%

* = Outside of Control Limits.

Matrix Spike/Matrix Spike Duplicate Summary

Page 1 of 1

Job Number: D41448

Account: XTOKWR XTO Energy

Project: PCU 296-5A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
D41448-1MS	3V21982.D	1	12/06/12	BD	n/a	n/a	V3V1284
D41448-1MSD	3V21983.D	1	12/06/12	BD	n/a	n/a	V3V1284
D41448-1	3V21984.D	1	12/06/12	BD	n/a	n/a	V3V1284

The QC reported here applies to the following samples:

Method: SW846 8260B

D41448-1

CAS No.	Compound	D41448-1		Spike	MS	MS	MSD	MSD	Limits	
		ug/kg	Q	ug/kg	ug/kg	%	ug/kg	%	RPD	Rec/RPD
71-43-2	Benzene	128		5590	5620	98	5970	104	6	64-139/30
100-41-4	Ethylbenzene	46.6	J	5590	5520	98	5860	104	6	68-136/30
108-88-3	Toluene	327		5590	5210	87	5460	92	5	60-130/30
1330-20-7	Xylene (total)	399	J	16800	16600	97	17600	103	6	58-142/30

CAS No.	Surrogate Recoveries	MS	MSD	D41448-1	Limits
2037-26-5	Toluene-D8	95%	95%	98%	64-130%
460-00-4	4-Bromofluorobenzene	96%	96%	94%	62-131%
17060-07-0	1,2-Dichloroethane-D4	112%	115%	111%	70-130%

* = Outside of Control Limits.

6.3.1
6



GC/MS Volatiles

Raw Data

7

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\V3120612.S\
 Data File : 3V21984.D
 Acq On : 6 Dec 2012 3:20 pm
 Operator : BRETD
 Sample : D41448-1
 Misc : MS5060,V3V1284,5.015,,100,5,1
 ALS Vial : 11 Sample Multiplier: 1

Quant Time: Dec 07 08:33:55 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	140932	50.00	ug/l	0.00
35) 1,4-Difluorobenzene	12.656	114	244735	50.00	ug/l	0.00
53) Chlorobenzene-d5	15.296	117	274569	50.00	ug/l	0.00
74) 1,4-Dichlorobenzene-d4	17.281	152	149790	50.00	ug/l	0.00

System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	12.251	102	19031	55.63	ug/l	0.00
Spiked Amount 50.000	Range 70 - 130		Recovery	=	111.26%	
61) Toluene-d8	14.048	98	325346	49.15	ug/l	0.00
Spiked Amount 50.000	Range 70 - 130		Recovery	=	98.30%	
69) 4-Bromofluorobenzene	16.242	95	127254	46.78	ug/l	0.00
Spiked Amount 50.000	Range 70 - 130		Recovery	=	93.56%	

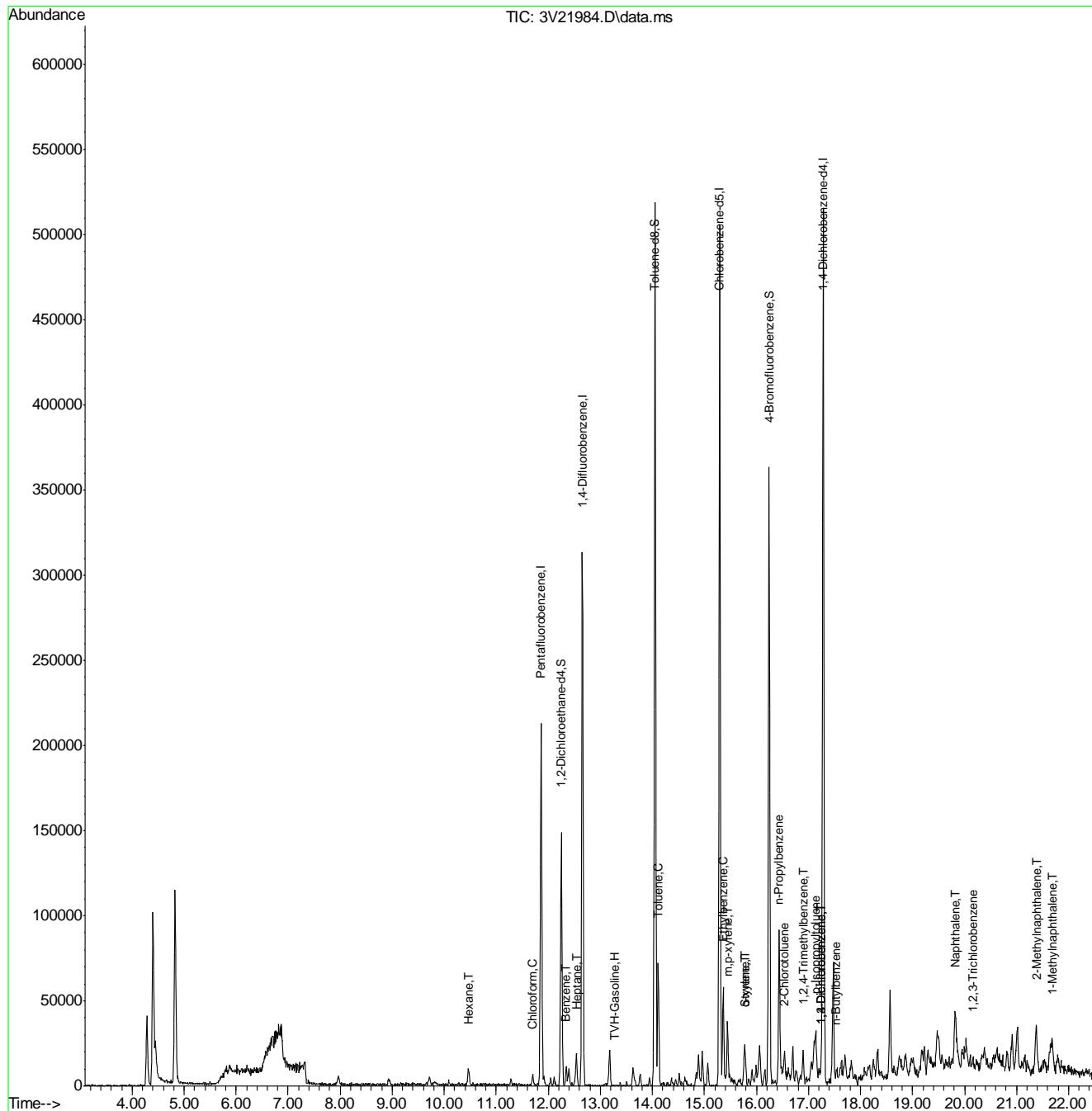
Target Compounds					Qvalue
1) TVH-Gasoline	13.285	TIC	888792m	187.86	ug/l
29) Chloroform	11.703	83	4615	1.29	ug/l 99
41) Hexane	10.471	57	4856	1.36	ug/l 100
43) Heptane	12.537	43	8730	1.94	ug/l 87
50) Benzene	12.354	78	9491	1.15	ug/l 100
62) Toluene	14.112	92	17551	2.93	ug/l 99
66) Ethylbenzene	15.360	91	4686	0.42	ug/l 99
71) Styrene	15.793	104	1924	0.45	ug/l 97
72) m,p-xylene	15.446	106	11659	2.54	ug/l 94
73) o-xylene	15.796	106	2337	1.03	ug/l 95
77) n-Propylbenzene	16.421	91	4777	0.38	ug/l # 89
78) 2-Chlorotoluene	16.534	91	503	0.06	ug/l 89
82) 1,2,4-Trimethylbenzene	16.893	105	8748	0.97	ug/l 91
84) 1,3-Dichlorobenzene	17.233	146	2080	0.40	ug/l # 85
85) 1,4-Dichlorobenzene	17.233	146	2079	0.38	ug/l 90
86) p-Isopropyltoluene	17.153	119	9040	0.95	ug/l # 86
88) n-Butylbenzene	17.538	91	4977	0.57	ug/l # 85
91) Naphthalene	19.834	128	22111	5.80	ug/l 100
93) 1,2,3-Trichlorobenzene	20.165	180	2704	0.94	ug/l 94
94) 2-Methylnaphthalene	21.381	142	18930	6.49	ug/l # 95
95) 1-Methylnaphthalene	21.689	142	13979	5.22	ug/l # 93

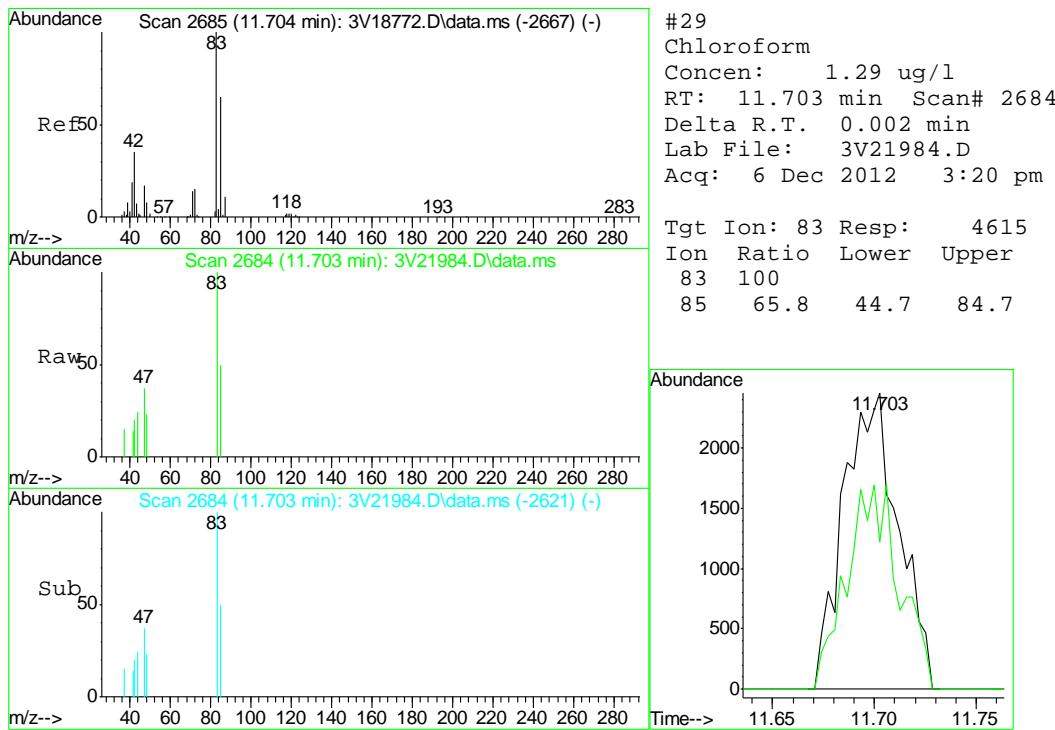
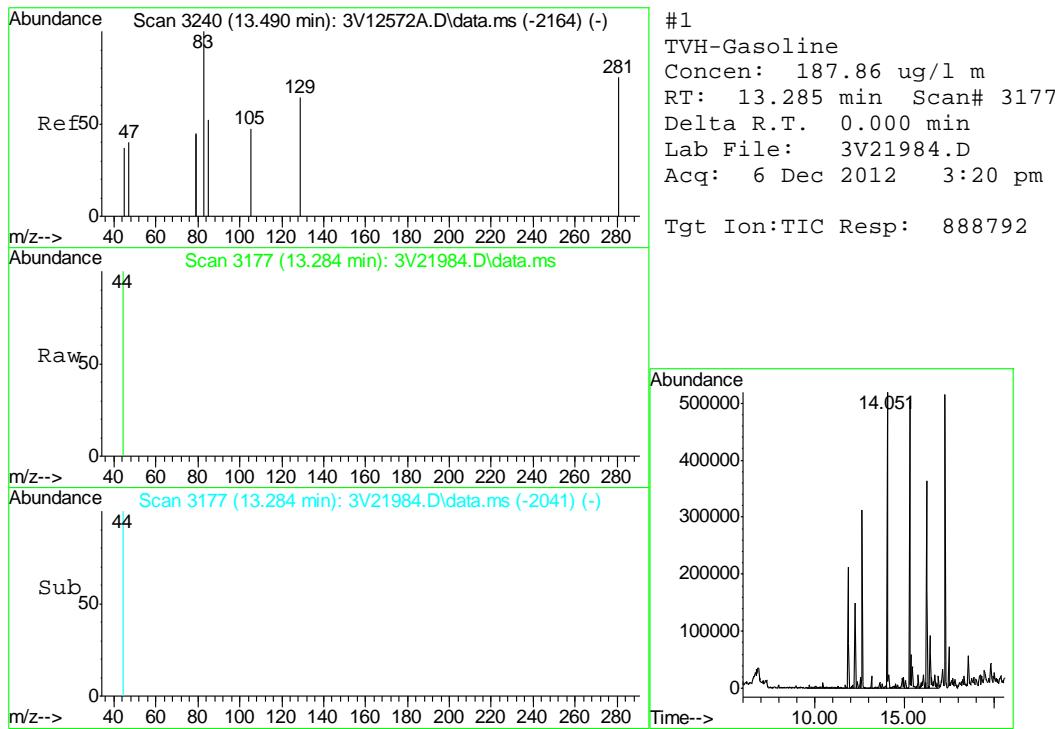
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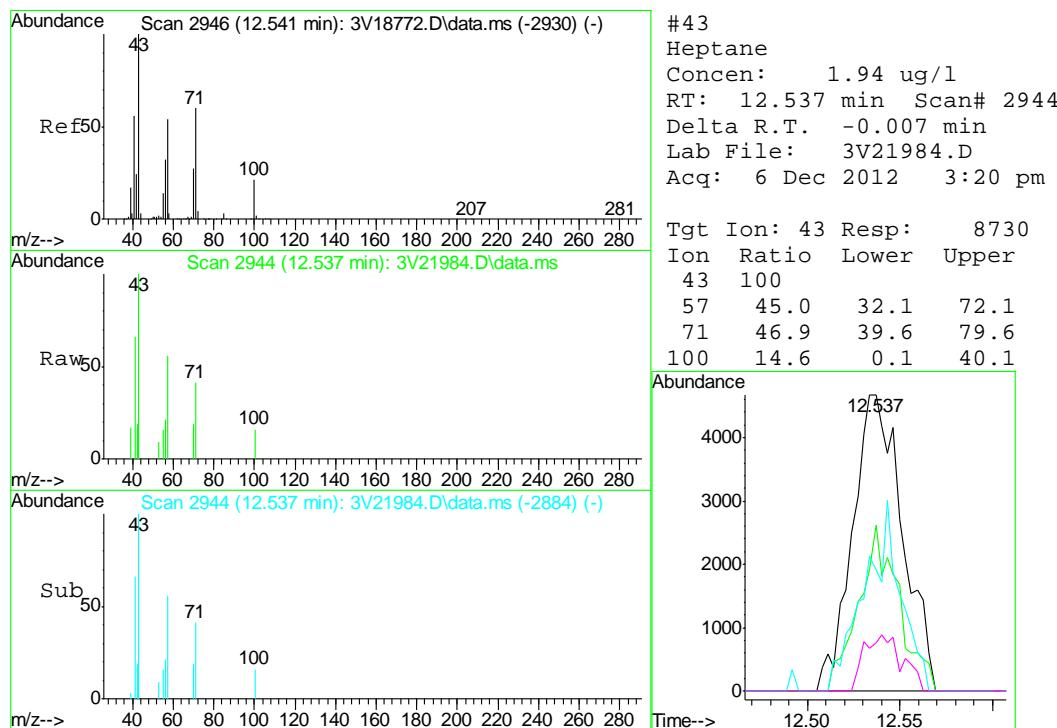
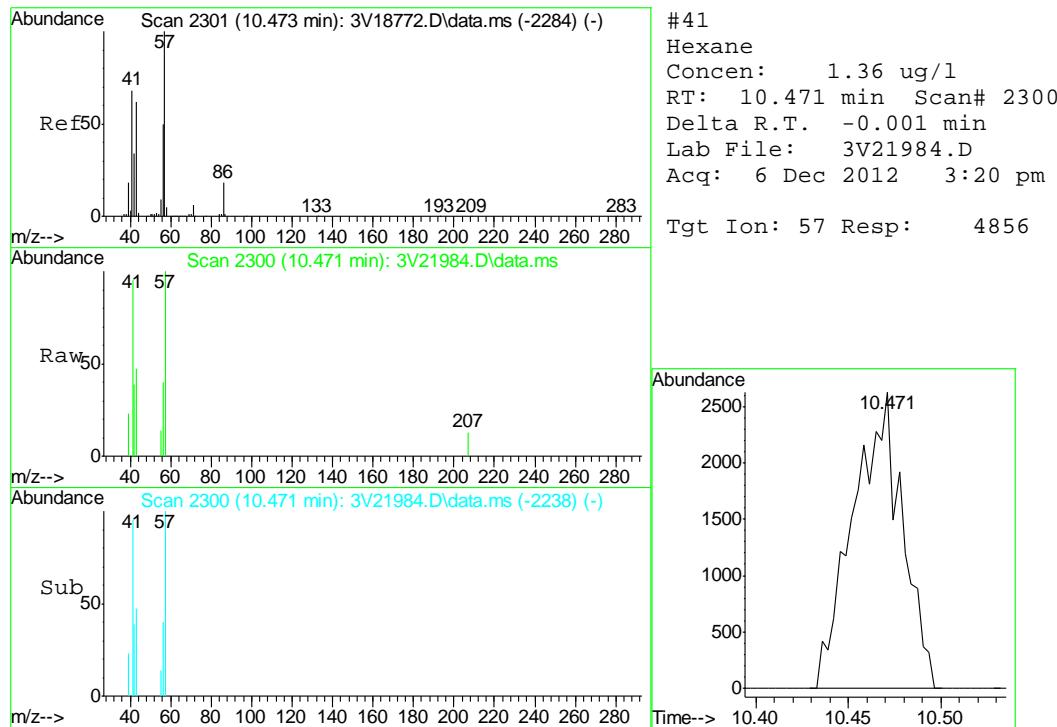
Quantitation Report (QT Reviewed)

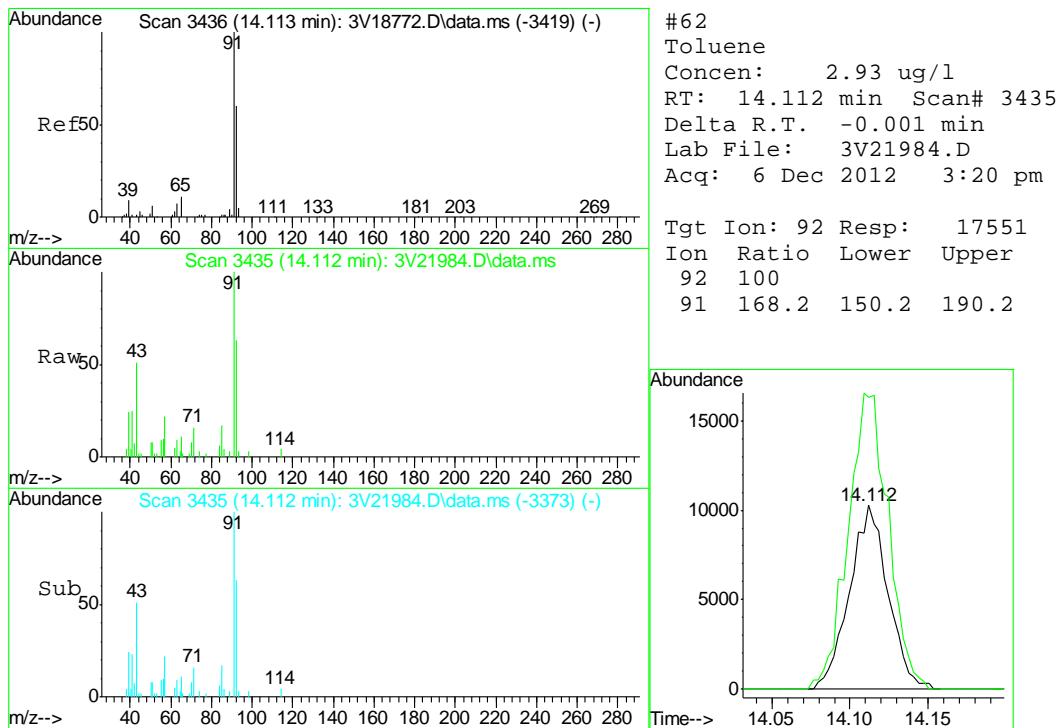
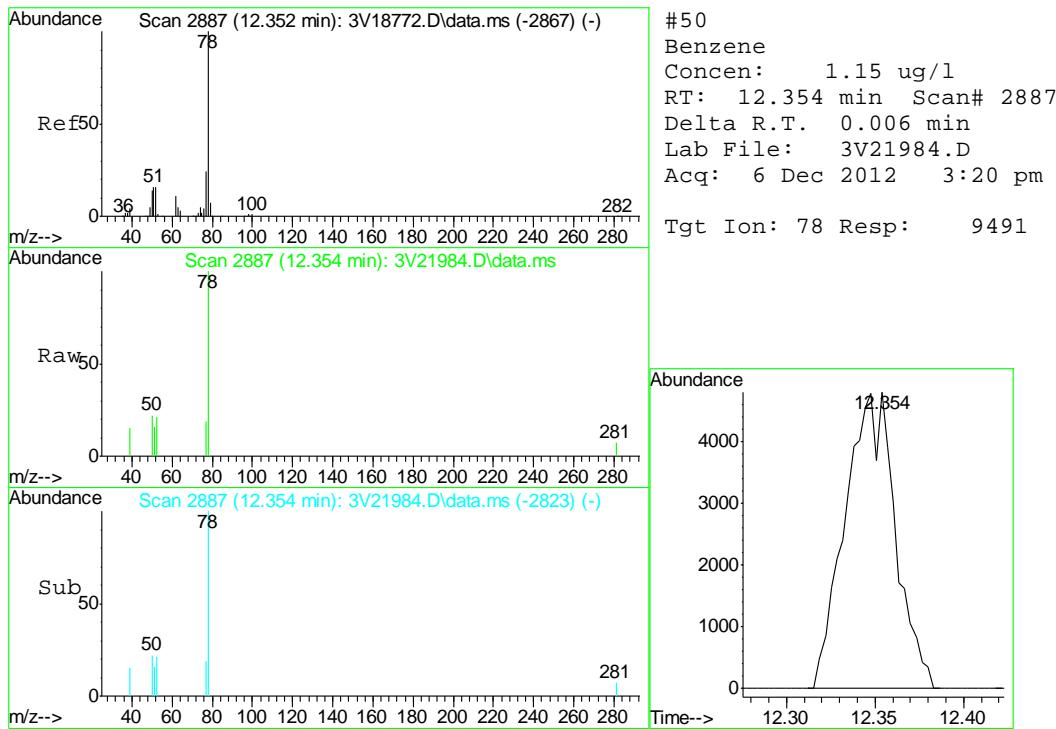
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 Operator : BRETD
 Sample : D41448-1
 Misc : MS5060,V3V1284,5.015,,100,5,1
 ALS Vial : 11 Sample Multiplier: 1

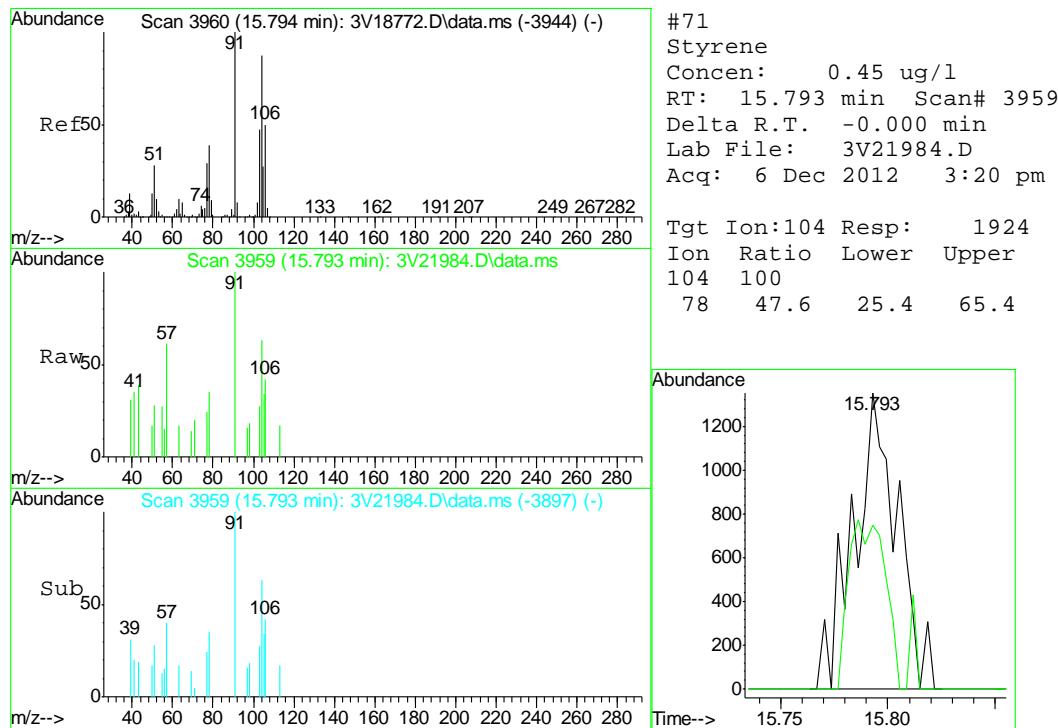
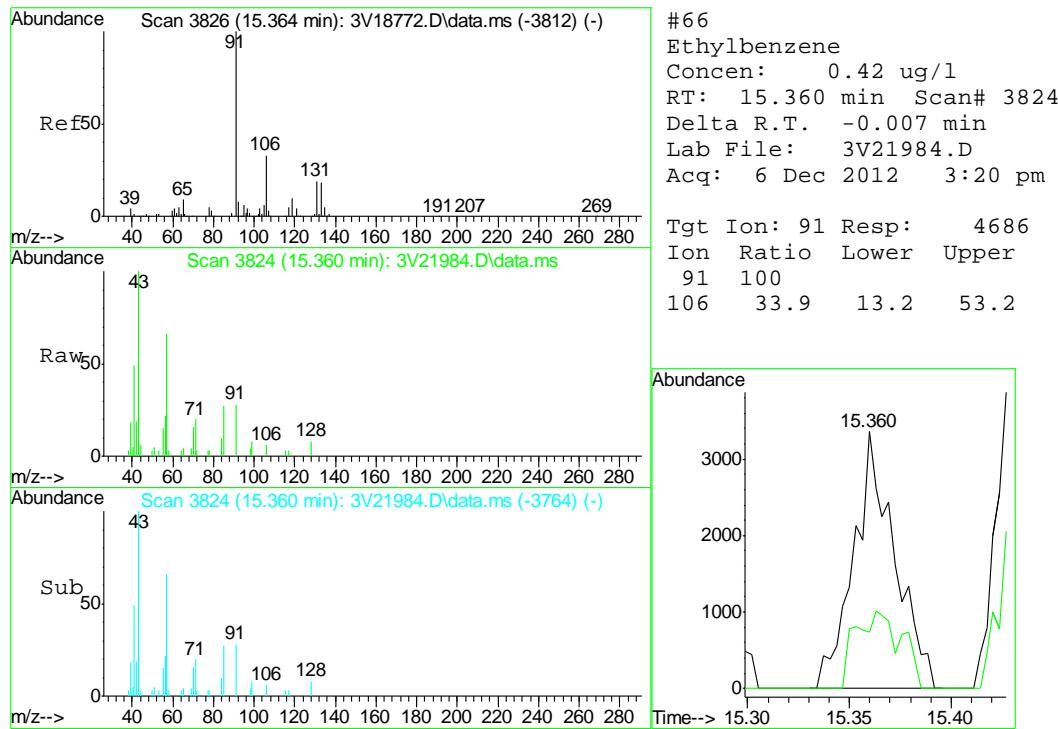
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 Response via : Initial Calibration

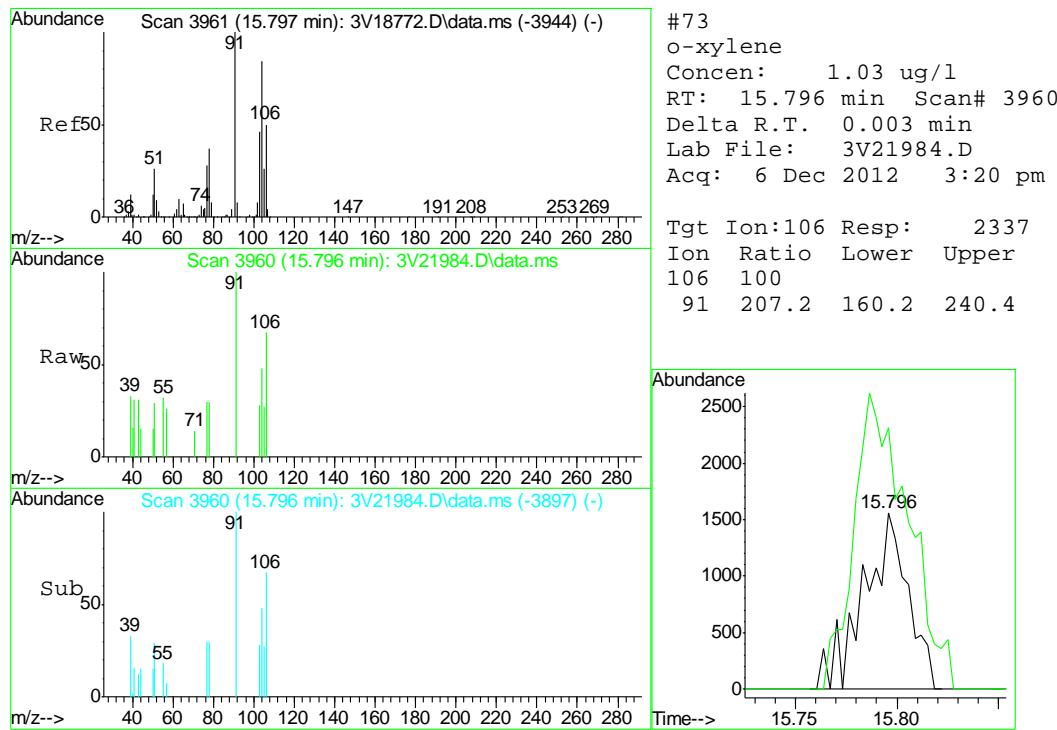
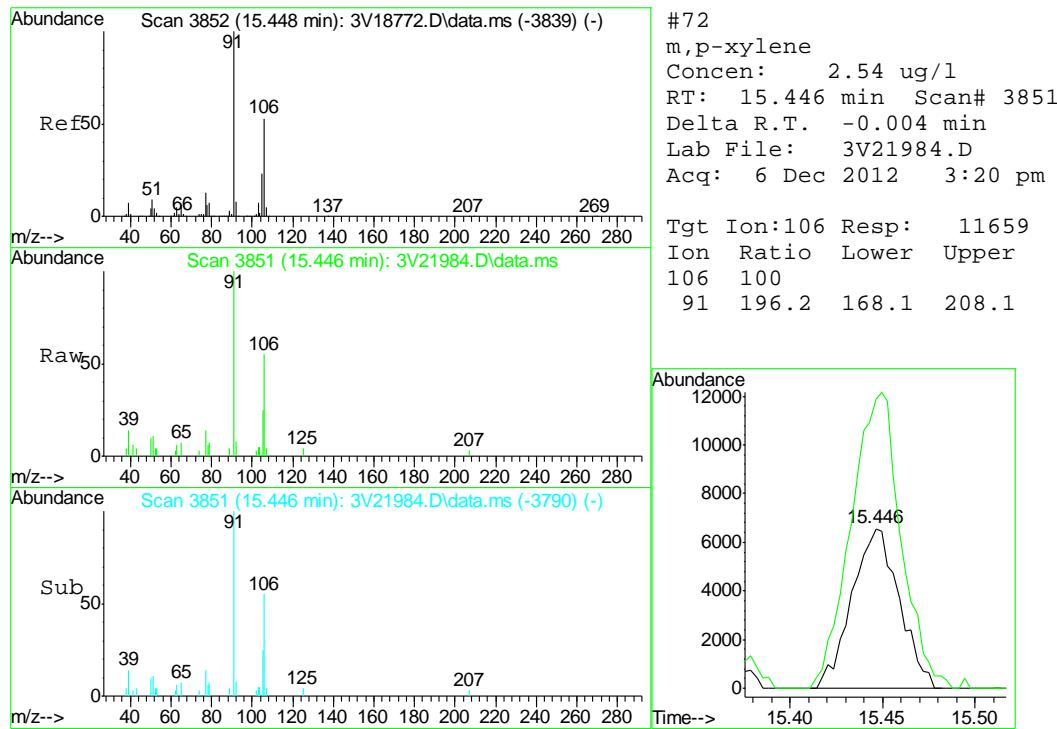


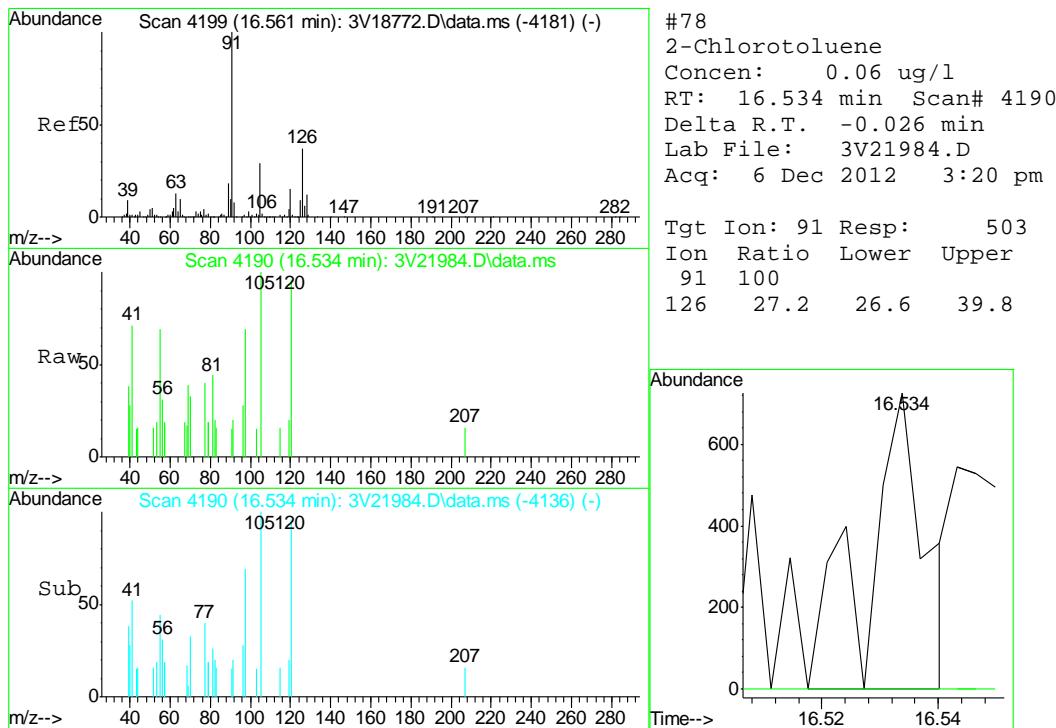
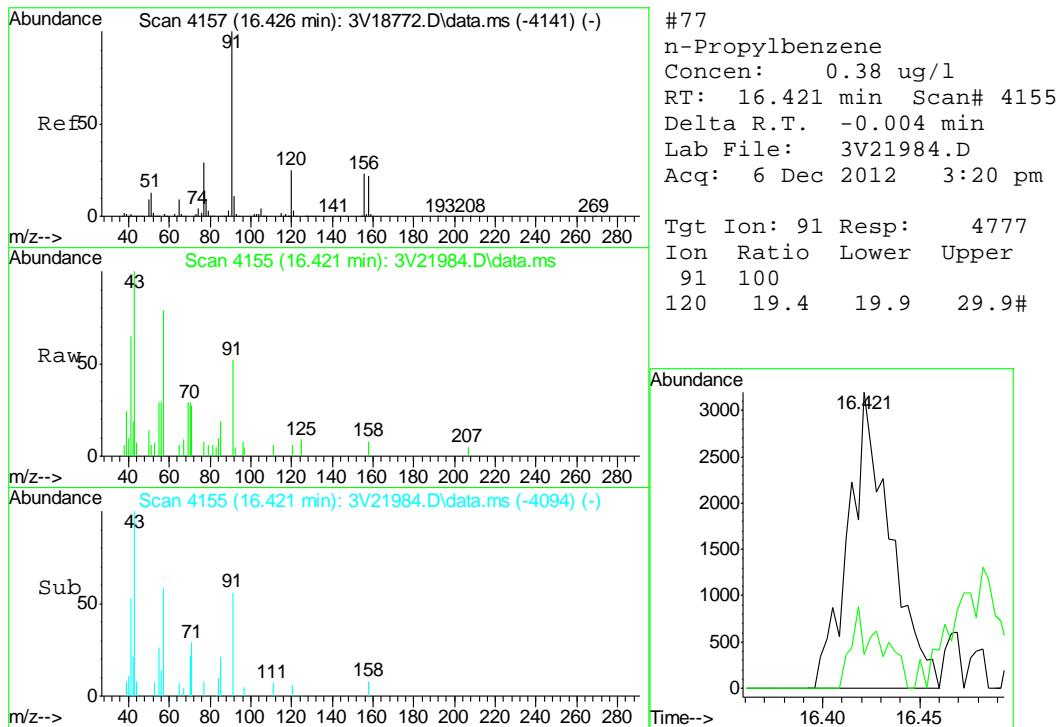


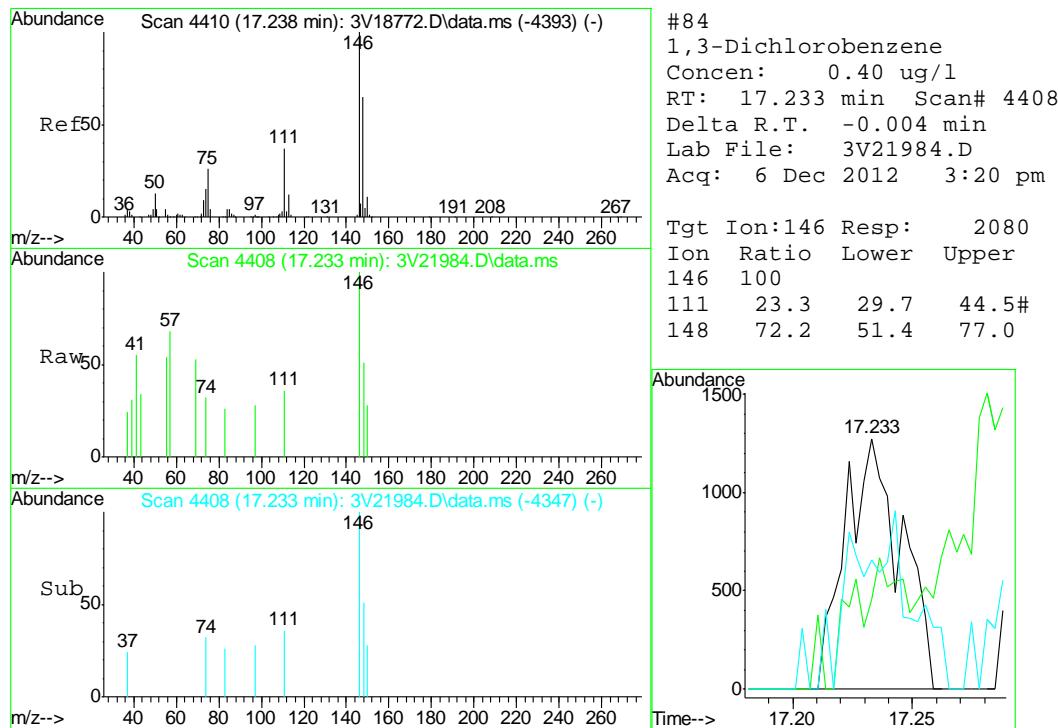
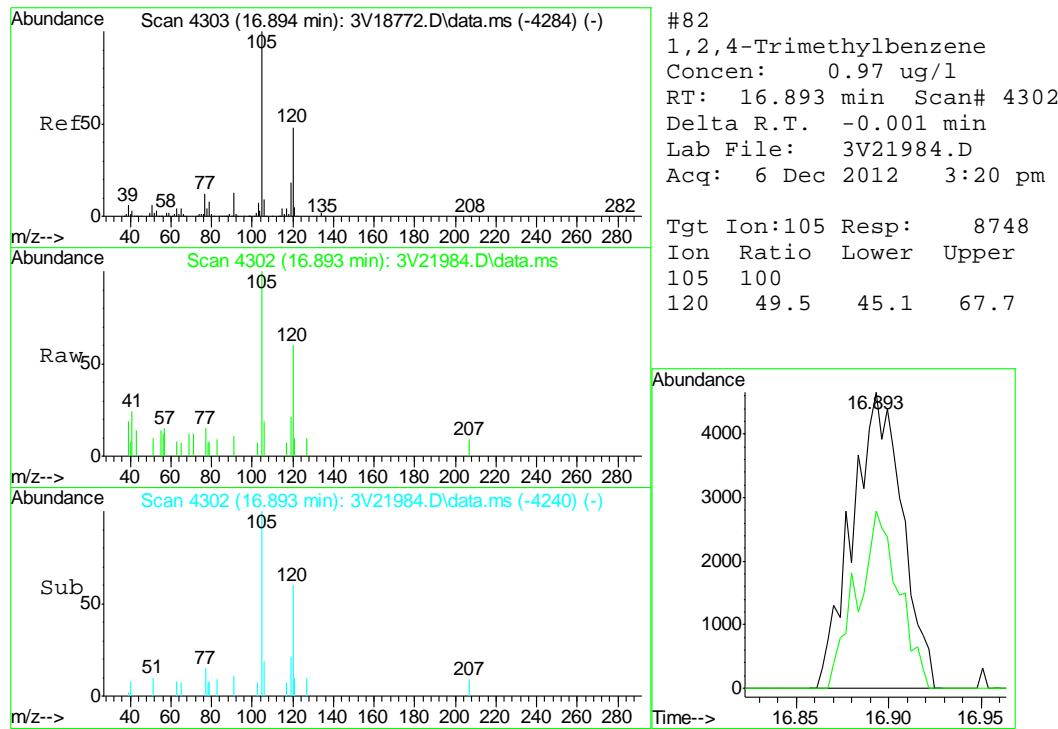


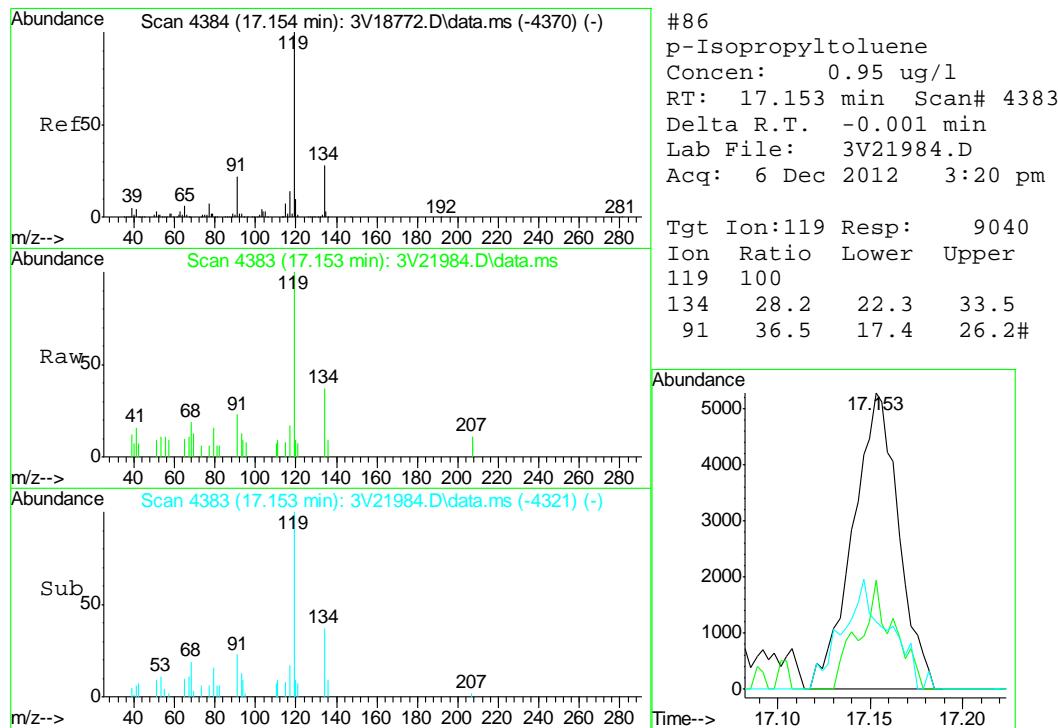
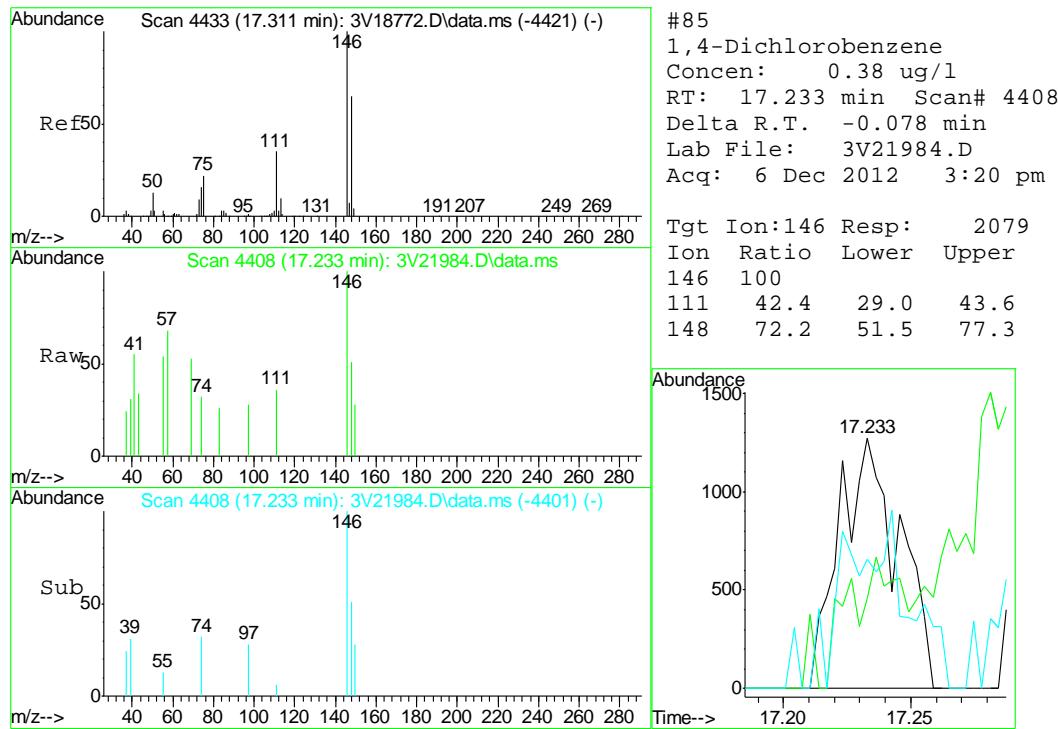


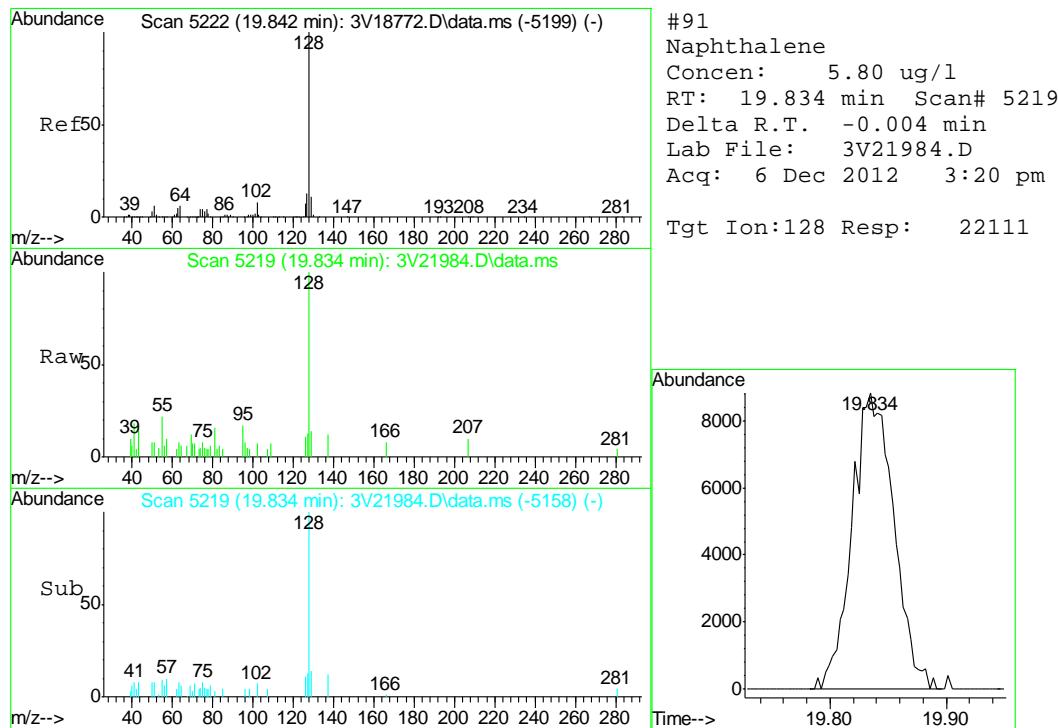
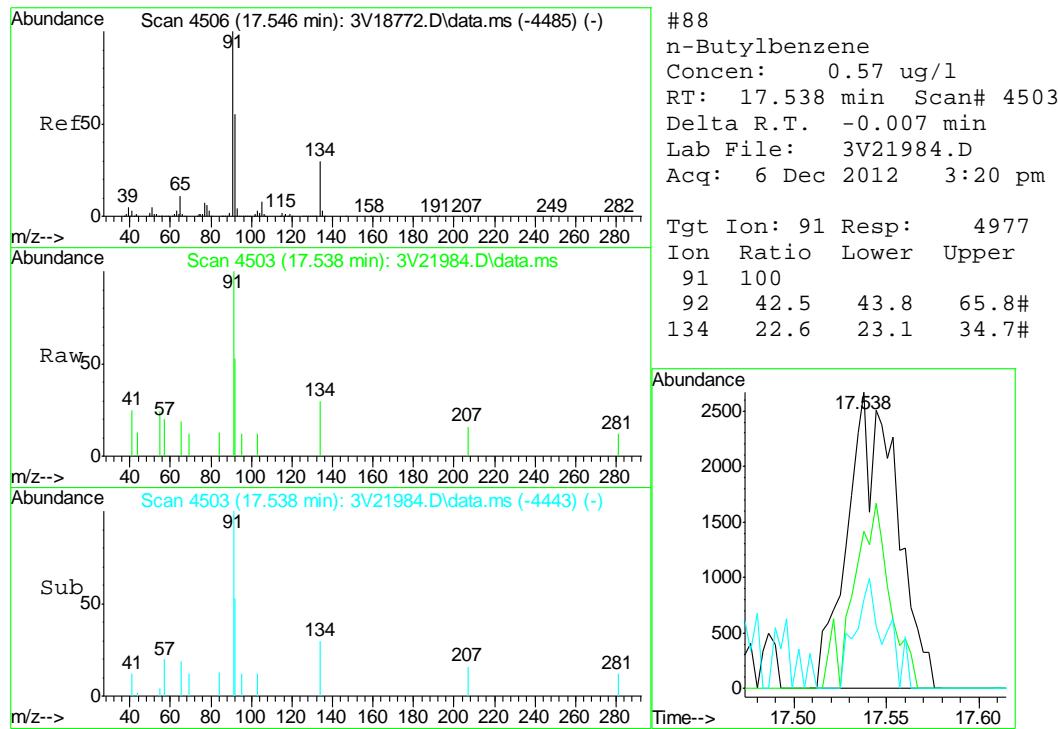


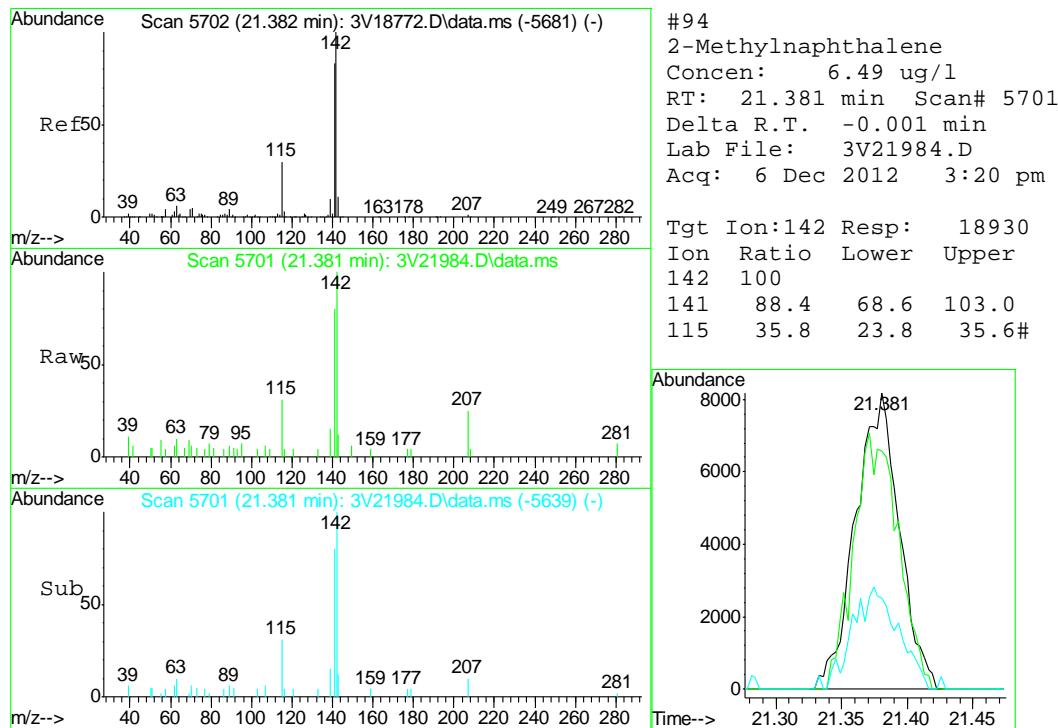
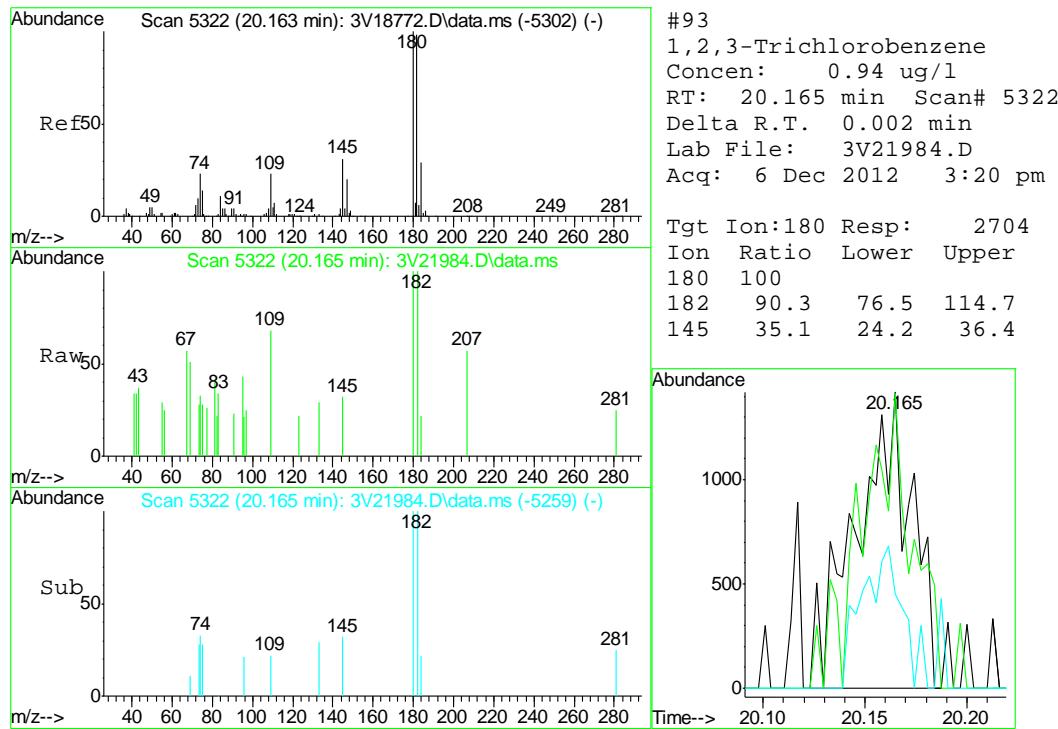


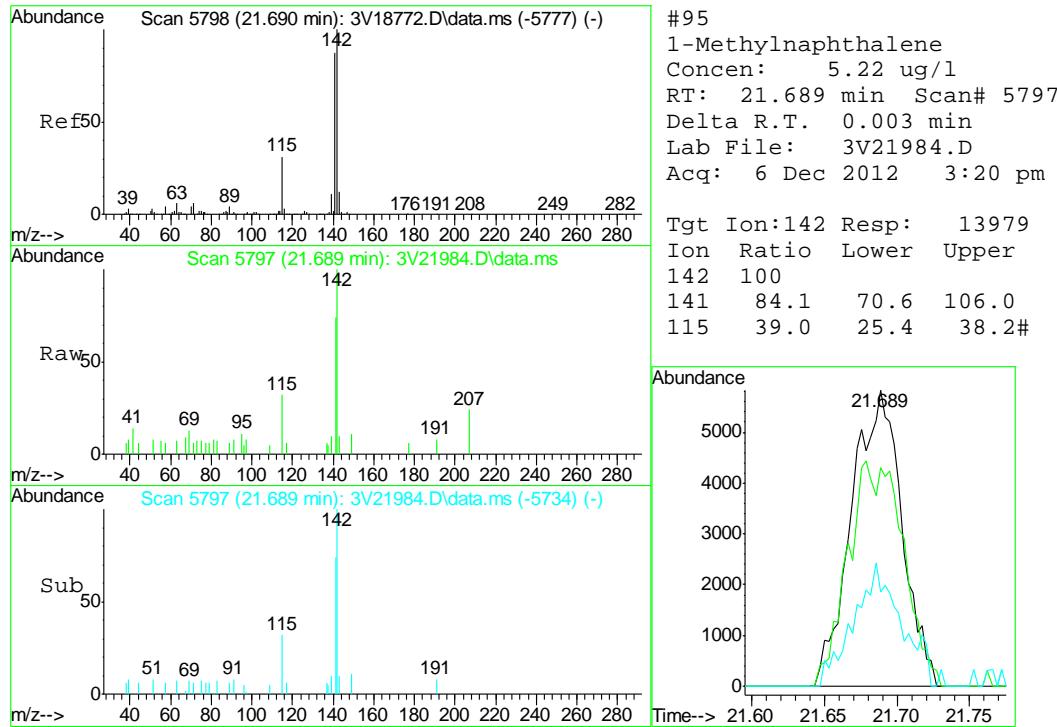












Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\V3120612.S\
 Data File : 3V21977.D
 Acq On : 6 Dec 2012 11:41 am
 Operator : BRETD
 Sample : MB
 Misc : MS5060,V3V1284,5.00,,100,5,1
 ALS Vial : 4 Sample Multiplier: 1

Quant Time: Dec 07 08:18:07 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	11.858	168	134900	50.00	ug/l	0.00
35) 1,4-Difluorobenzene	12.653	114	239166	50.00	ug/l	0.00
53) Chlorobenzene-d5	15.290	117	253145	50.00	ug/l	0.00
74) 1,4-Dichlorobenzene-d4	17.279	152	133239	50.00	ug/l	0.00

System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	12.249	102	19278	58.87	ug/l	0.00
Spiked Amount	50.000	Range	70 - 130	Recovery	=	117.74%
61) Toluene-d8	14.045	98	314722	51.57	ug/l	0.00
Spiked Amount	50.000	Range	70 - 130	Recovery	=	103.14%
69) 4-Bromofluorobenzene	16.239	95	108246	43.16	ug/l	0.00
Spiked Amount	50.000	Range	70 - 130	Recovery	=	86.32%

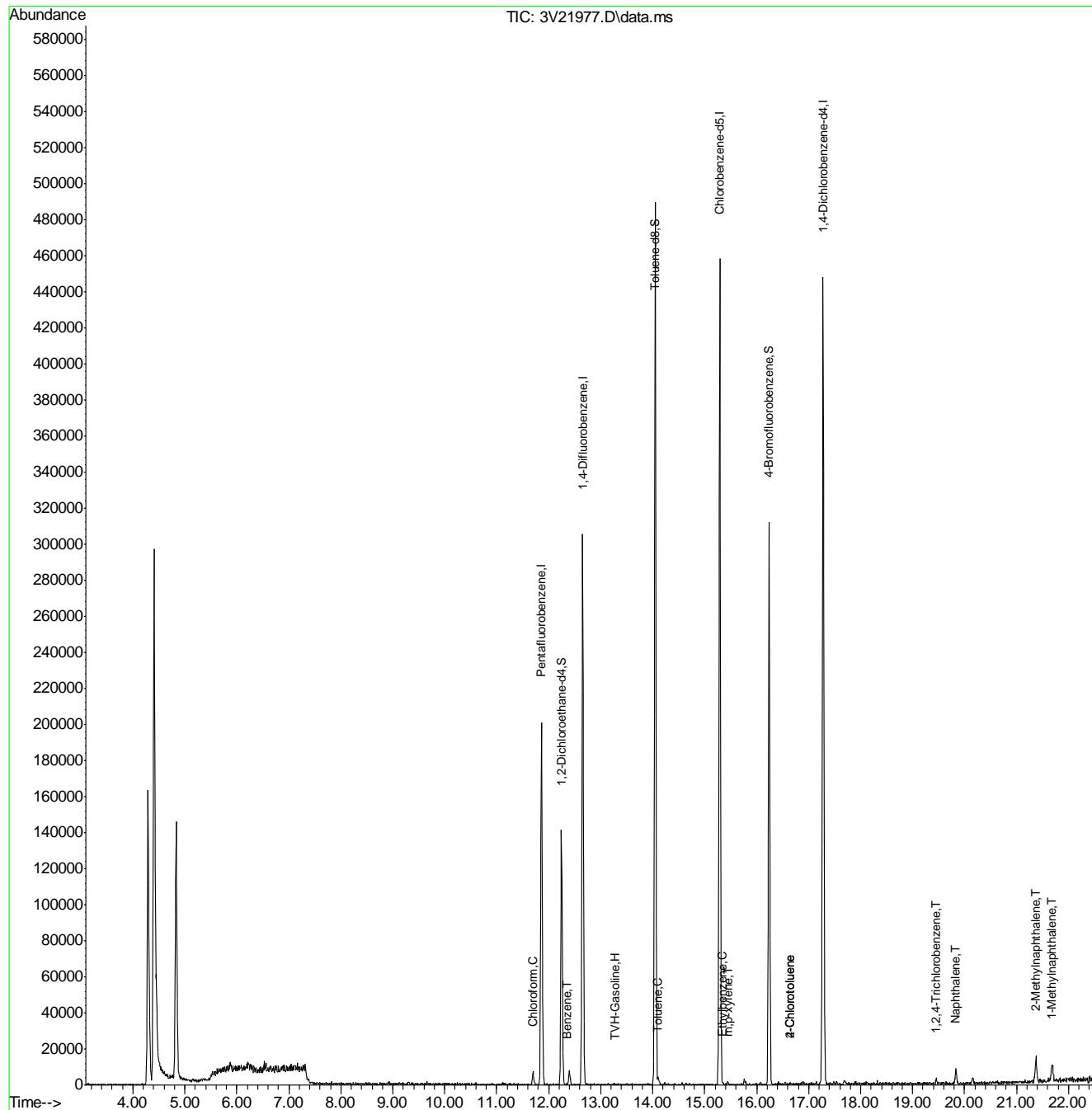
Target Compounds					Qvalue
1) TVH-Gasoline	13.285	TIC	36710m	160.24	ug/l
29) Chloroform	11.697	83	4791	1.40	ug/l
50) Benzene	12.345	78	178	0.02	ug/l
62) Toluene	14.106	92	1520	0.27	ug/l
66) Ethylbenzene	15.354	91	401	0.04	ug/l
72) m,p-xylene	15.457	106	548	0.13	ug/l #
78) 2-Chlorotoluene	16.643	91	406	0.05	ug/l
79) 4-Chlorotoluene	16.643	91	406	0.06	ug/l
90) 1,2,4-Trichlorobenzene	19.453	180	1330	0.50	ug/l #
91) Naphthalene	19.832	128	11253	4.73	ug/l
94) 2-Methylnaphthalene	21.368	142	10507	4.05	ug/l
95) 1-Methylnaphthalene	21.673	142	7775	3.27	ug/l #

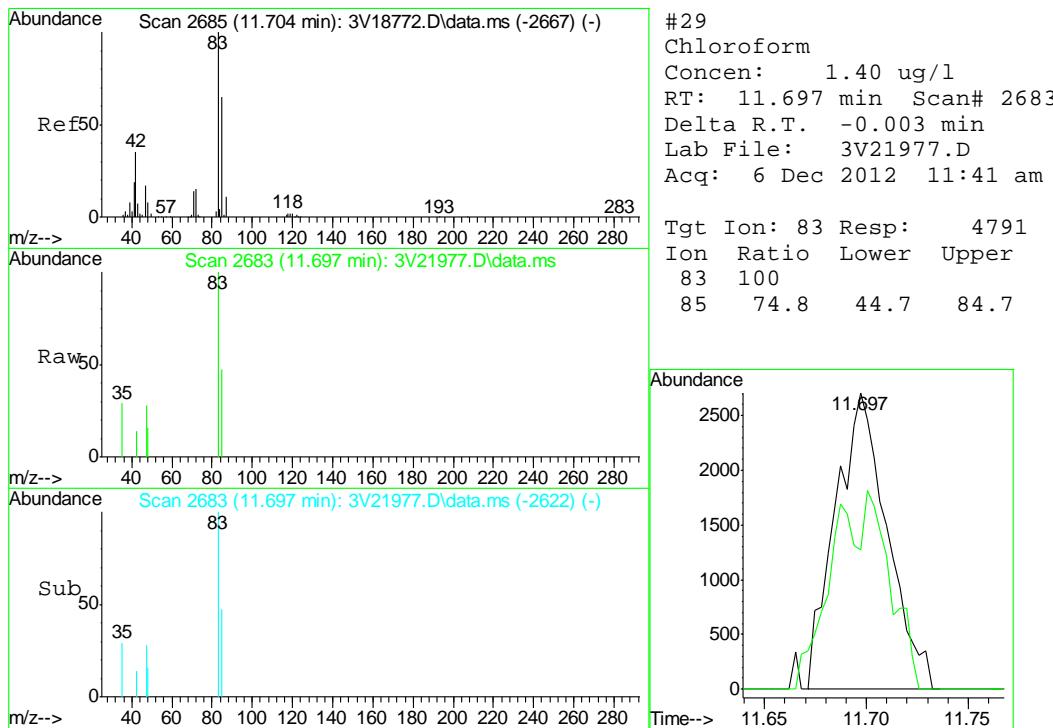
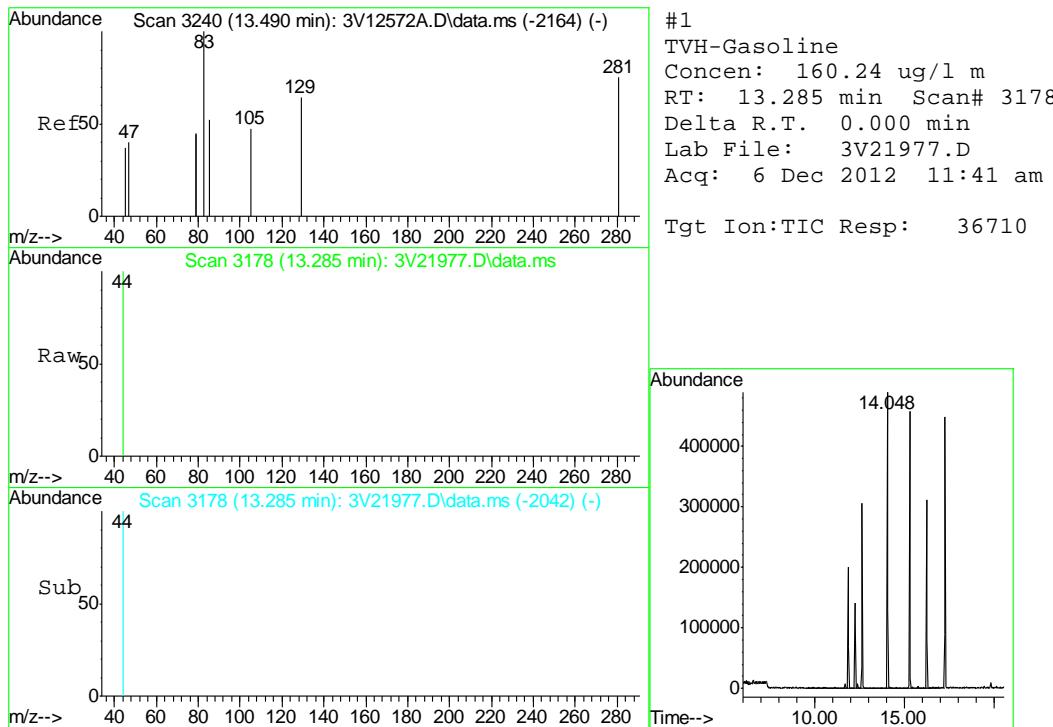
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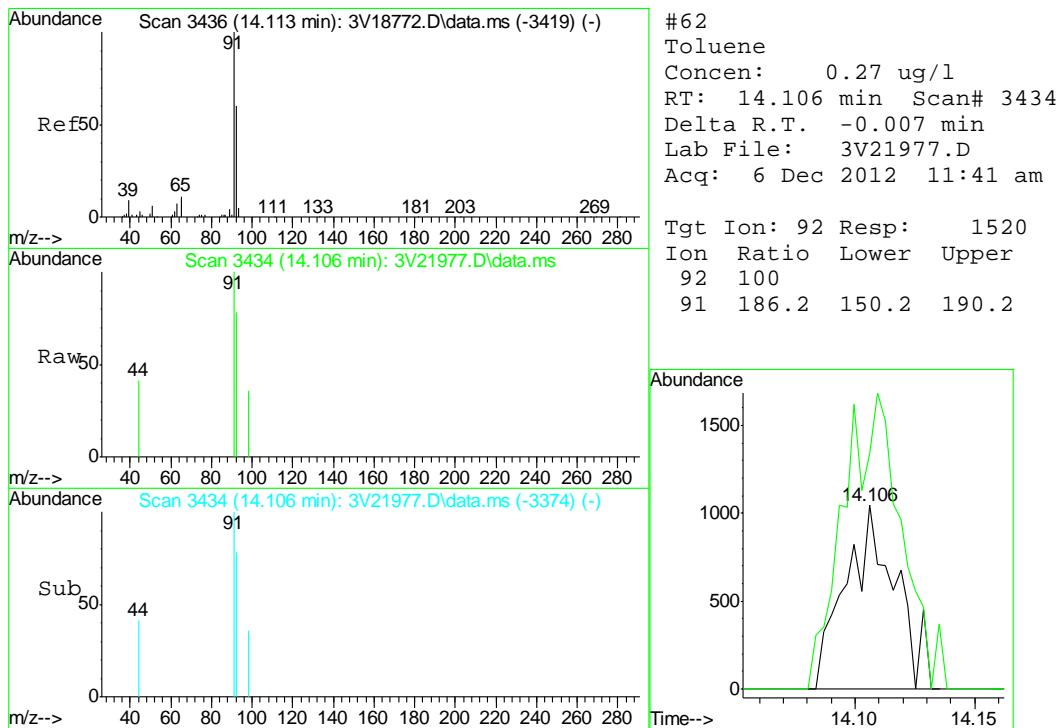
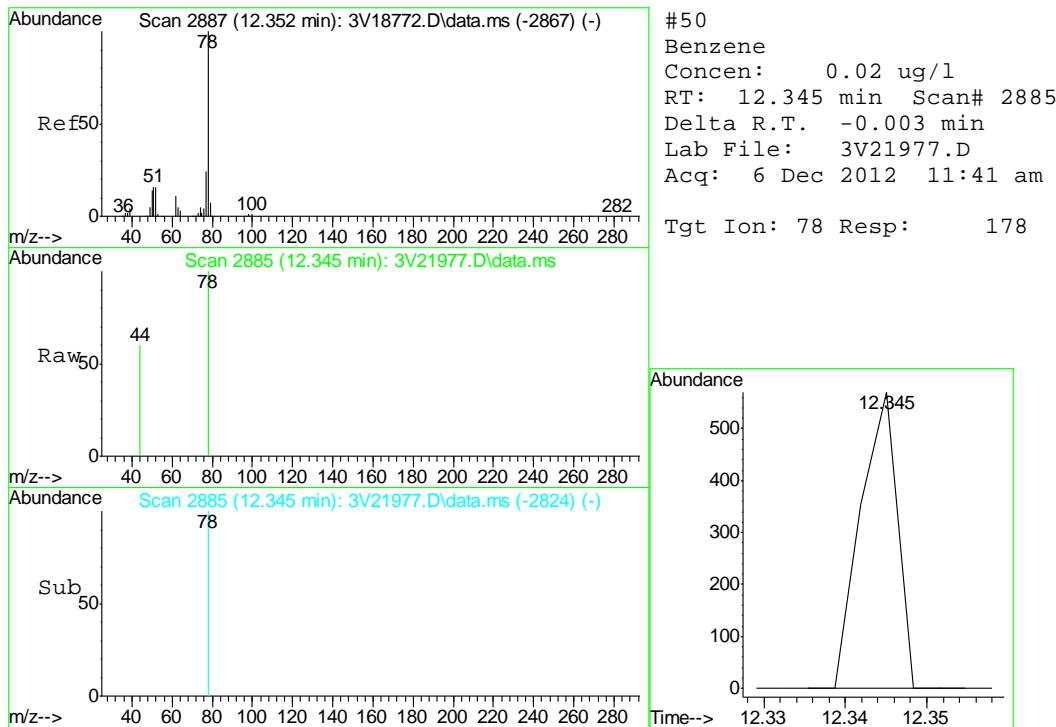
Quantitation Report (QT Reviewed)

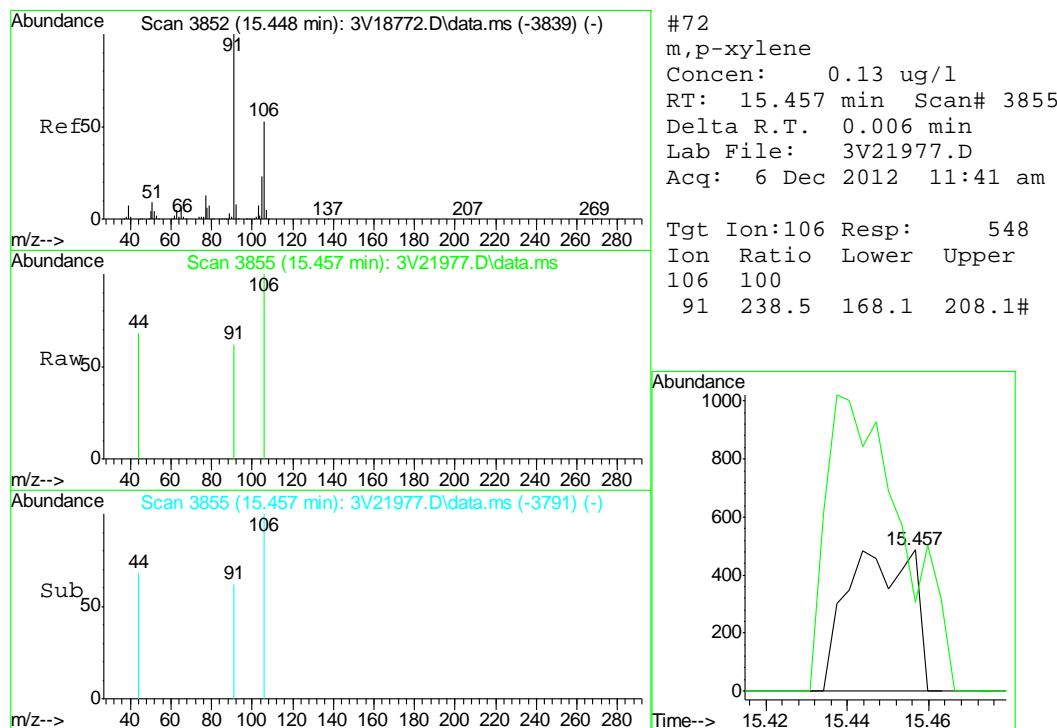
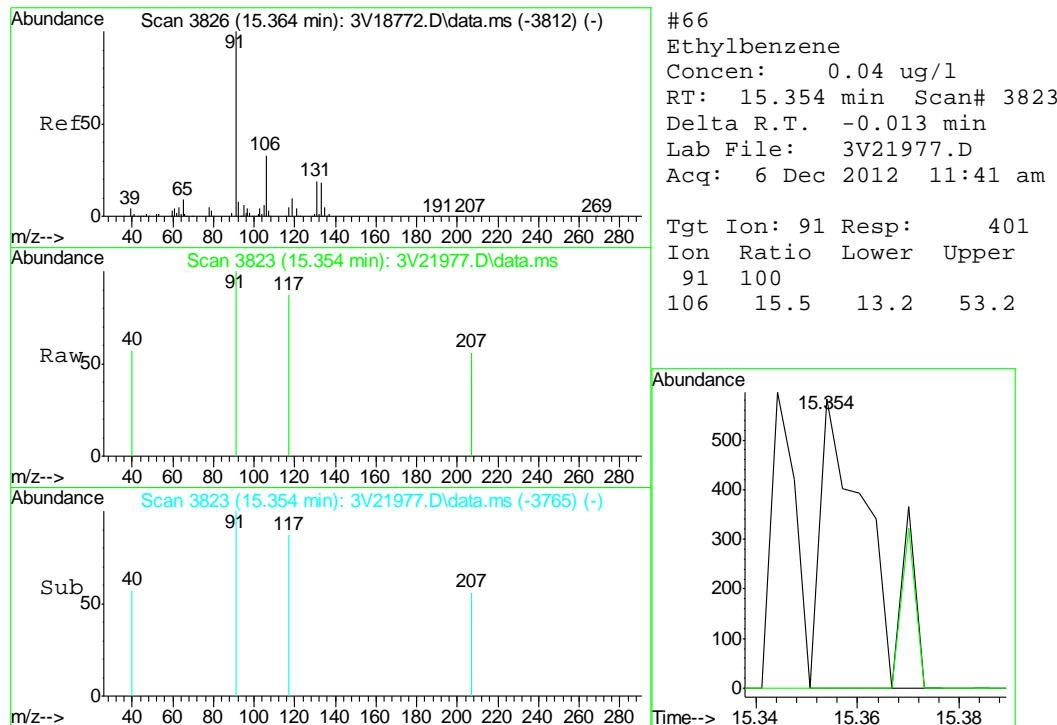
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 Data File : 3V21977.D
 Acq On : 6 Dec 2012 11:41 am
 Operator : BRETD
 Sample : MB
 Misc : MS5060,V3V1284,5.00,,100,5,1
 ALS Vial : 4 Sample Multiplier: 1

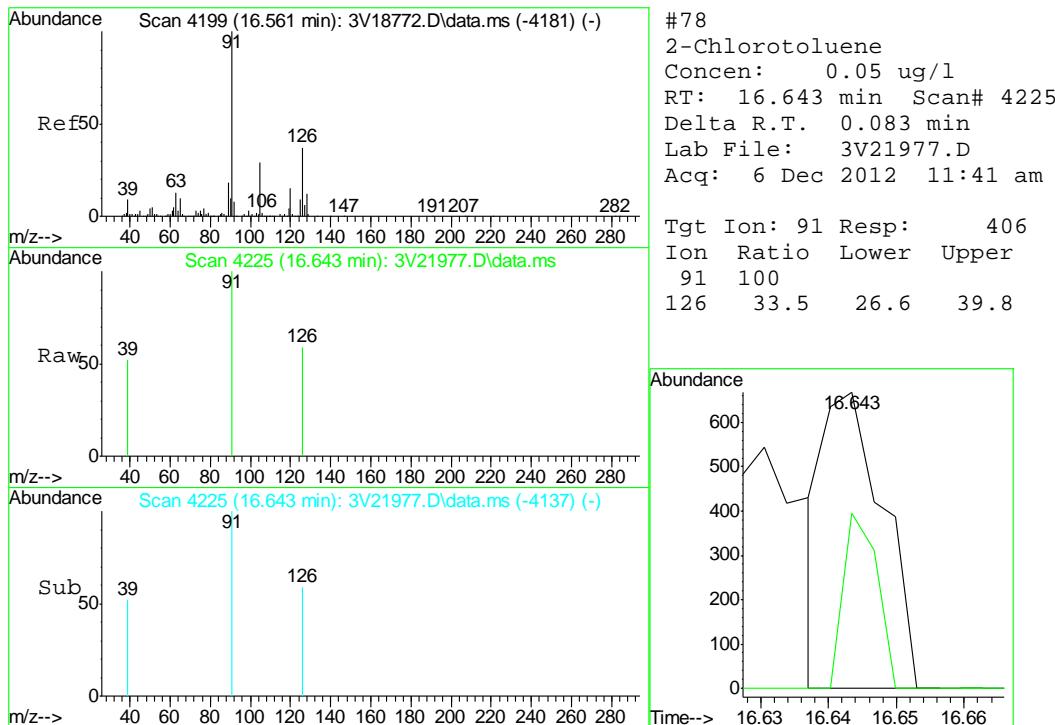
Quant Time: Dec 07 08:18:07 2012
 Quant Method : C:\msdchem\1\METHODS\V3AP1277TVH1277SOIL.M
 Quant Title : 8260
 QLast Update : Wed Nov 28 14:20:19 2012
 Response via : Initial Calibration



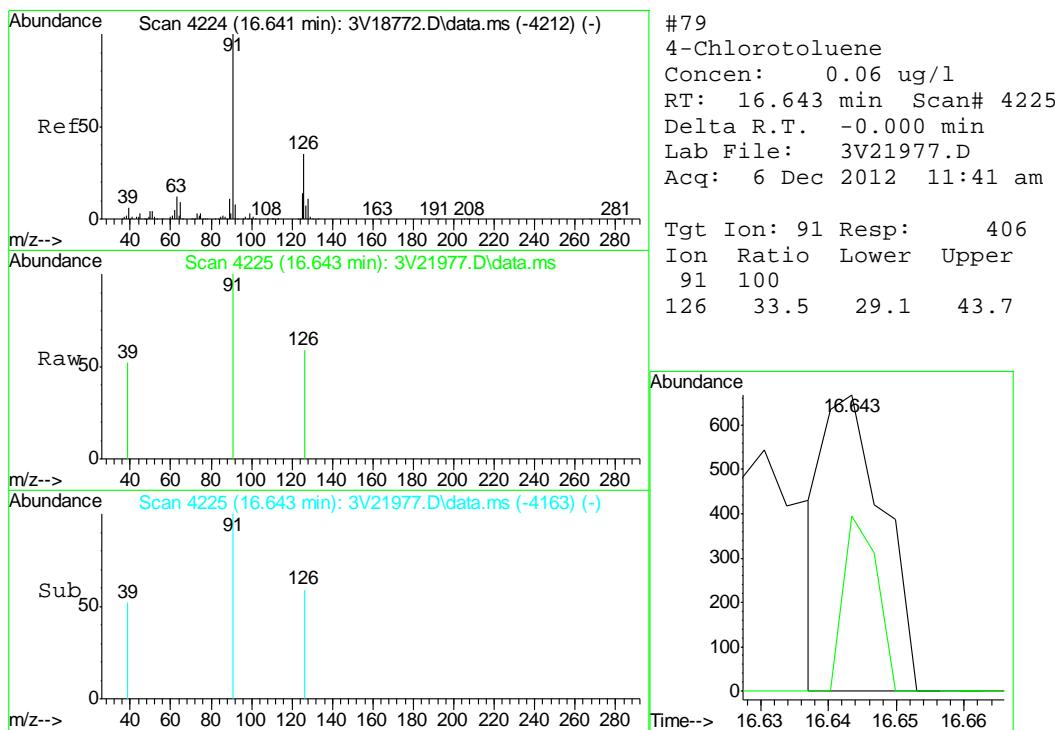


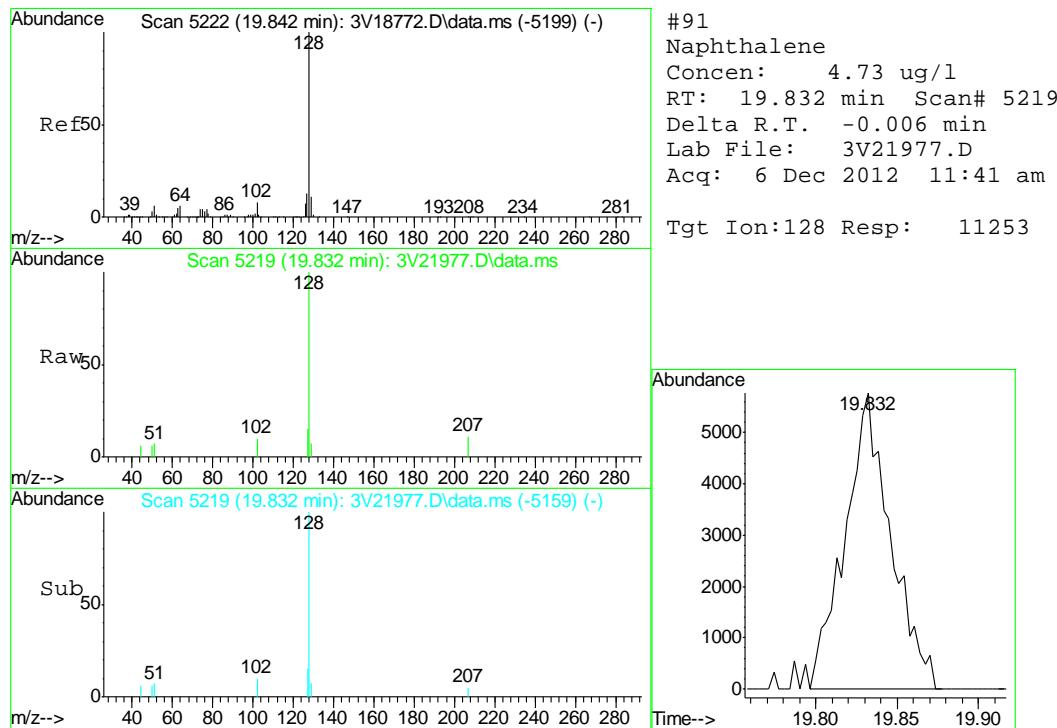
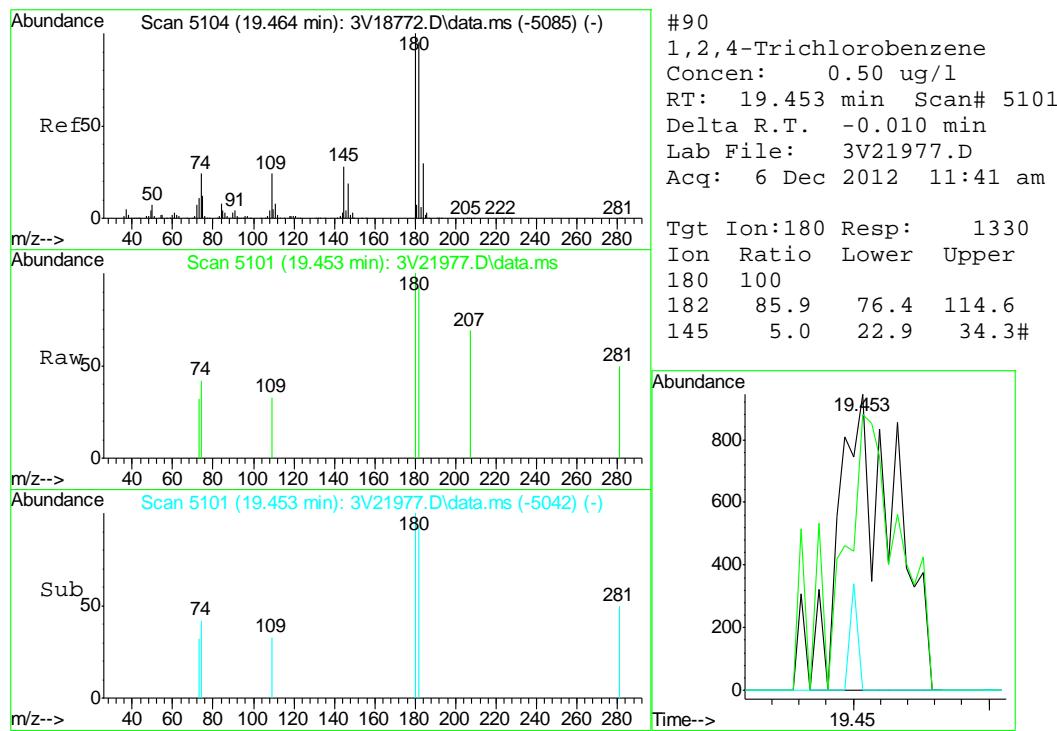


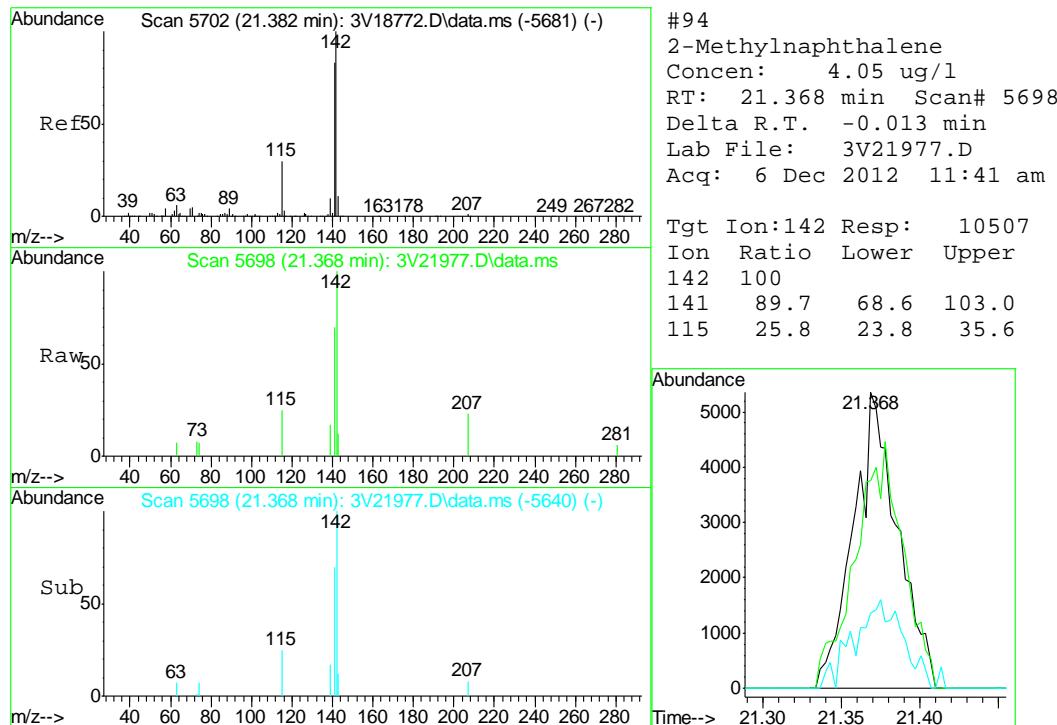




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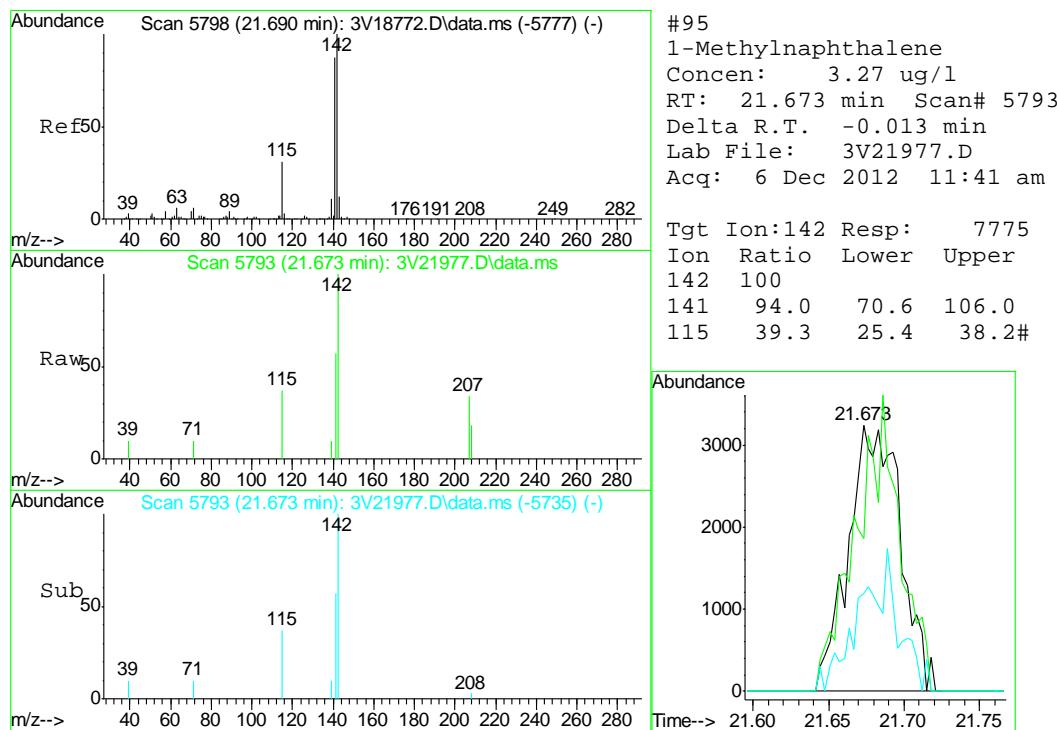






7.2.1

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GC/MS Semi-volatiles

QC Data Summaries

∞

Includes the following where applicable:

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

Method Blank Summary

Page 1 of 1

Job Number: D41448

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

D41448-1

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%	10-159%
321-60-8	2-Fluorobiphenyl	79%	19-131%
1718-51-0	Terphenyl-d14	91%	18-150%

Blank Spike Summary

Page 1 of 1

Job Number: D41448

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

D41448-1

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: D41448

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

D41448-1

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

Raw Data

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\121012\
 Data File : 3g12515.D
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 Operator : DONC
 Sample : D41448-1
 Misc : OP7075,E3G593,30.02,,,1,1
 ALS Vial : 11 Sample Multiplier: 1

Quant Time: Dec 10 14:58:19 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	5.671	136	139161	4.0000	ug/mL	-0.01
6) Acenaphthene-d10	7.385	164	90213	4.0000	ug/mL	0.00
15) Phenanthrene-d10	8.859	188	152658	4.0000	ug/mL	-0.02
19) Chrysene-d12	11.496	240	102798	4.0000	ug/mL	-0.02
24) Perylene-d12	12.873	264	83929	4.0000	ug/mL	-0.02

System Monitoring Compounds						
2) Nitrobenzene-d5	4.985	82	367269	26.4018	ug/mL	-0.01
Spiked Amount	50.000	Range	25 - 135	Recovery	=	52.80%
7) 2-Fluorobiphenyl	6.723	172	1027104	24.9801	ug/mL	0.00
Spiked Amount	50.000	Range	25 - 135	Recovery	=	49.96%
21) Terphenyl-d14	10.450	244	523060	34.5676	ug/mL	-0.02
Spiked Amount	50.000	Range	25 - 135	Recovery	=	69.14%

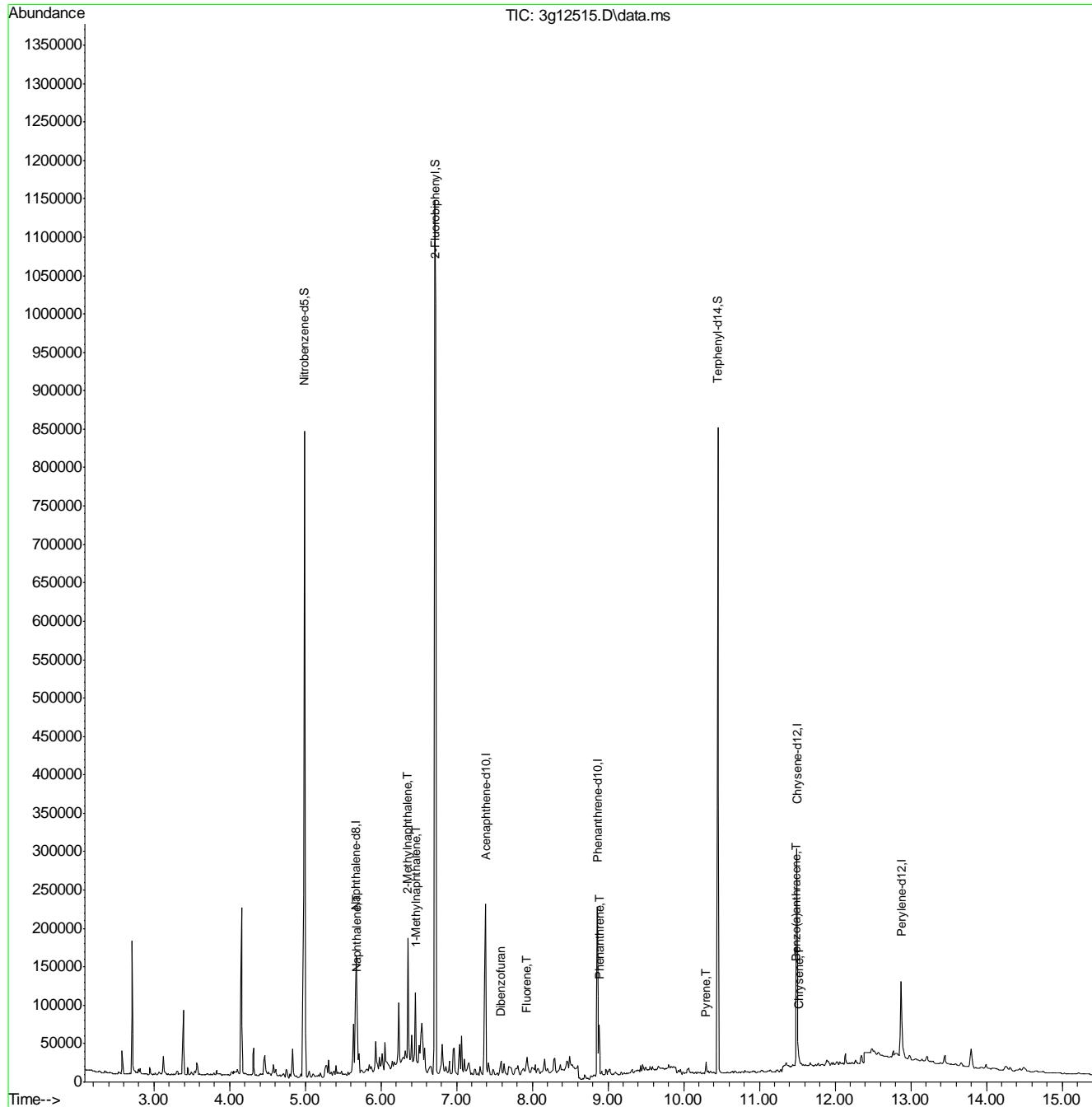
Target Compounds					Qvalue
3) N-Nitrosodimethylamine	0.000	74	0	N.D.	d
4) N-Nitrosodi-propylamine	0.000	70	0	N.D.	d
5) Naphthalene	5.683	128	44924	1.0670	ug/mL
8) 2-Methylnaphthalene	6.356	142	63460	1.9679	ug/mL
9) 1-Methylnaphthalene	6.456	142	34385	1.0759	ug/mL
10) Acenaphthylene	7.243	152	1350	N.D.	
11) Acenaphthene	7.408	154	778	Below Cal	# 1
12) Dibenzofuran	7.586	168	7529	0.1443	ug/mL
13) Fluorene	7.928	166	11166	0.2762	ug/mL#
14) Diphenylamine	0.000	169	0	N.D.	d
16) Phenanthrene	8.883	178	39297	0.6274	ug/mL
17) Anthracene	0.000	178	0	N.D.	d
18) Fluoranthene	0.000	202	0	N.D.	d
20) Pyrene	10.292	202	8564	0.1484	ug/mL#
22) Benzo(a)anthracene	11.483	228	3742	0.0758	ug/mL#
23) Chrysene	11.523	228	9657	0.1975	ug/mL#
25) Benzo(b)fluoranthene	0.000	252	0	N.D.	d
26) Benzo(k)fluoranthene	0.000	252	0	N.D.	d
27) Benzo(a)pyrene	0.000	252	0	N.D.	d
28) Indeno(1,2,3-cd)pyrene	14.093	276	1306	N.D.	
29) Dibenz(a,h)anthracene	14.103	278	1086	N.D.	
30) Benzo(g,h,i)perylene	14.093	276	1277	N.D.	

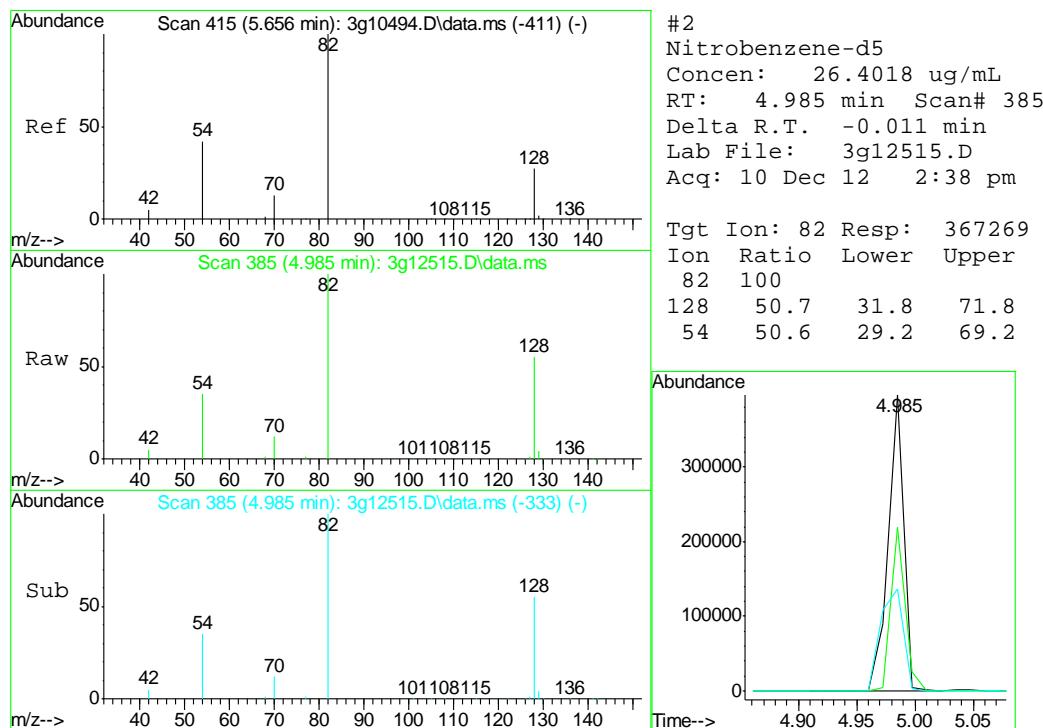
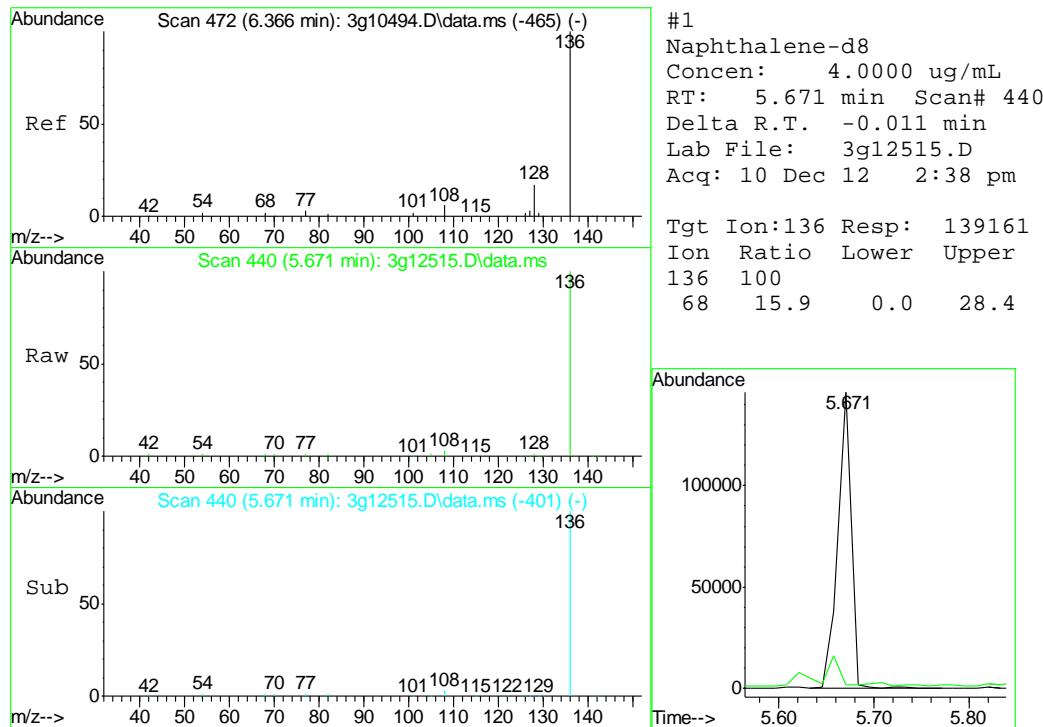
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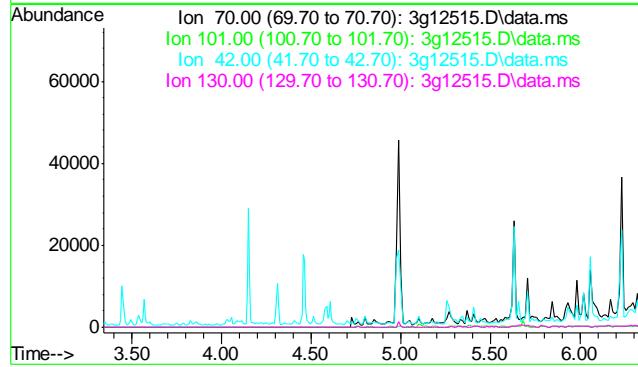
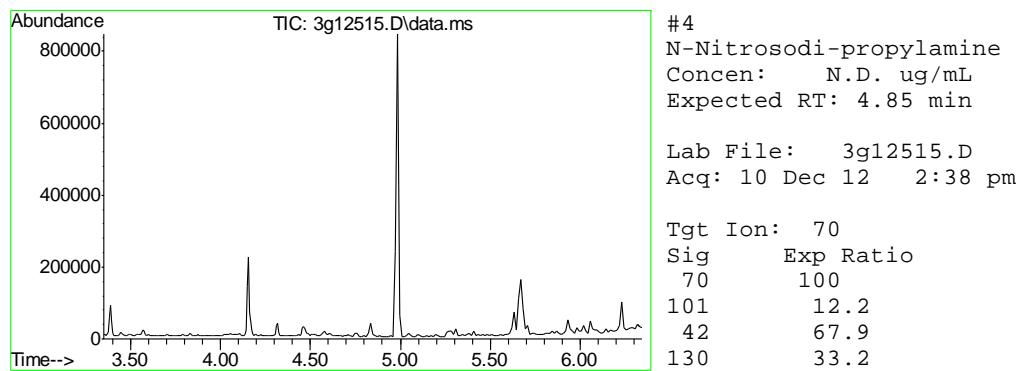
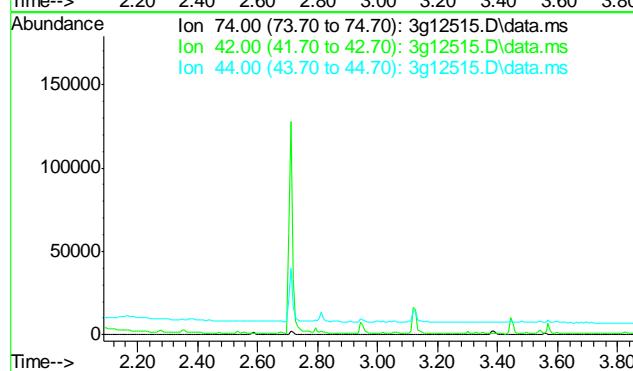
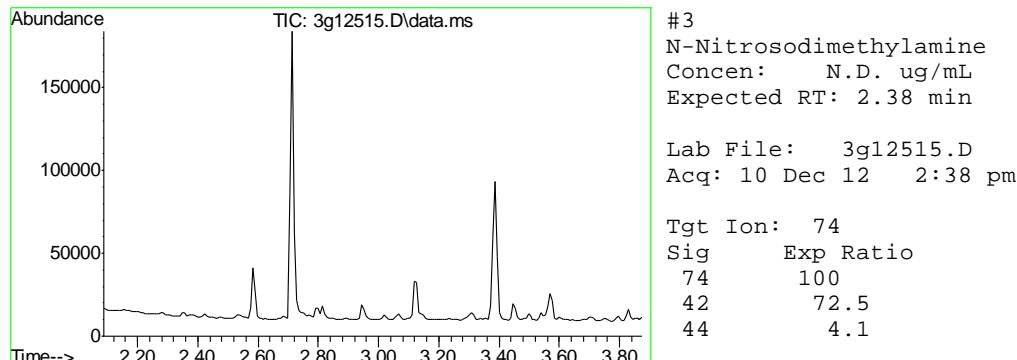
Quantitation Report (QT Reviewed)

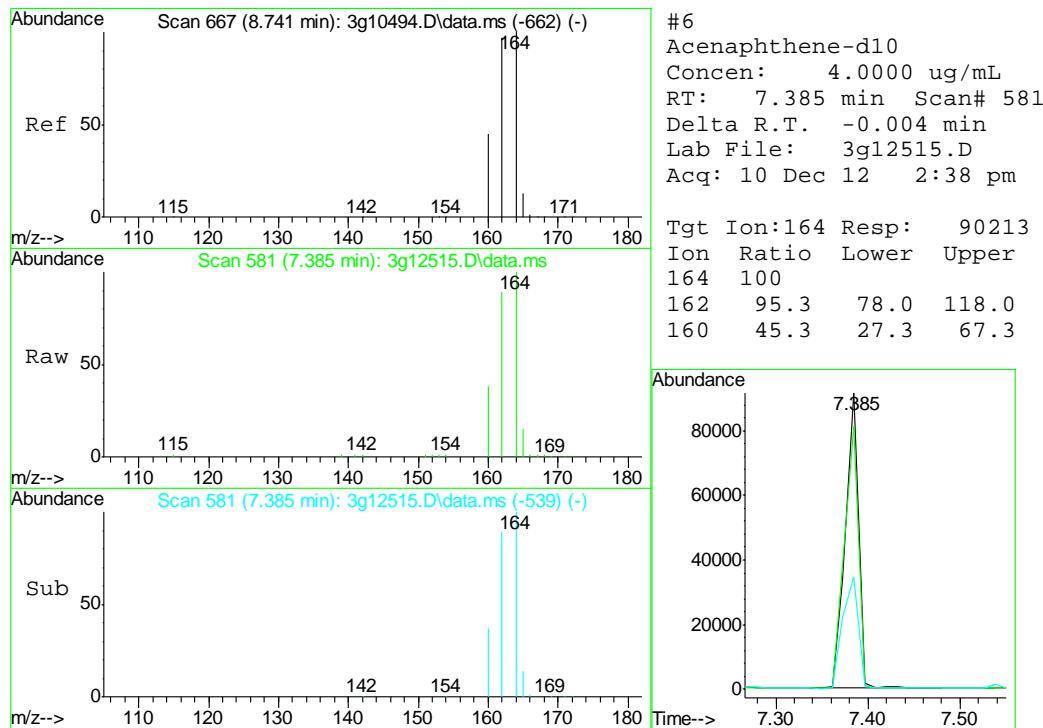
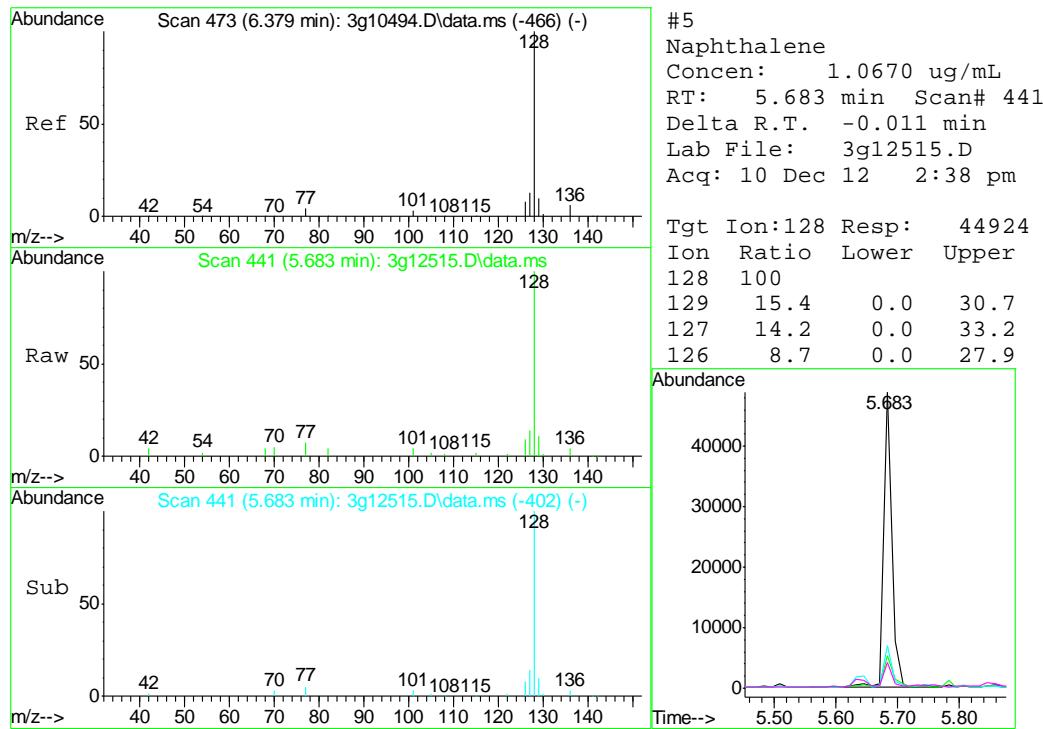
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 Operator : DONC
 Sample : D41448-1
 Misc : OP7075,E3G593,30.02,,,1,1
 ALS Vial : 11 Sample Multiplier: 1

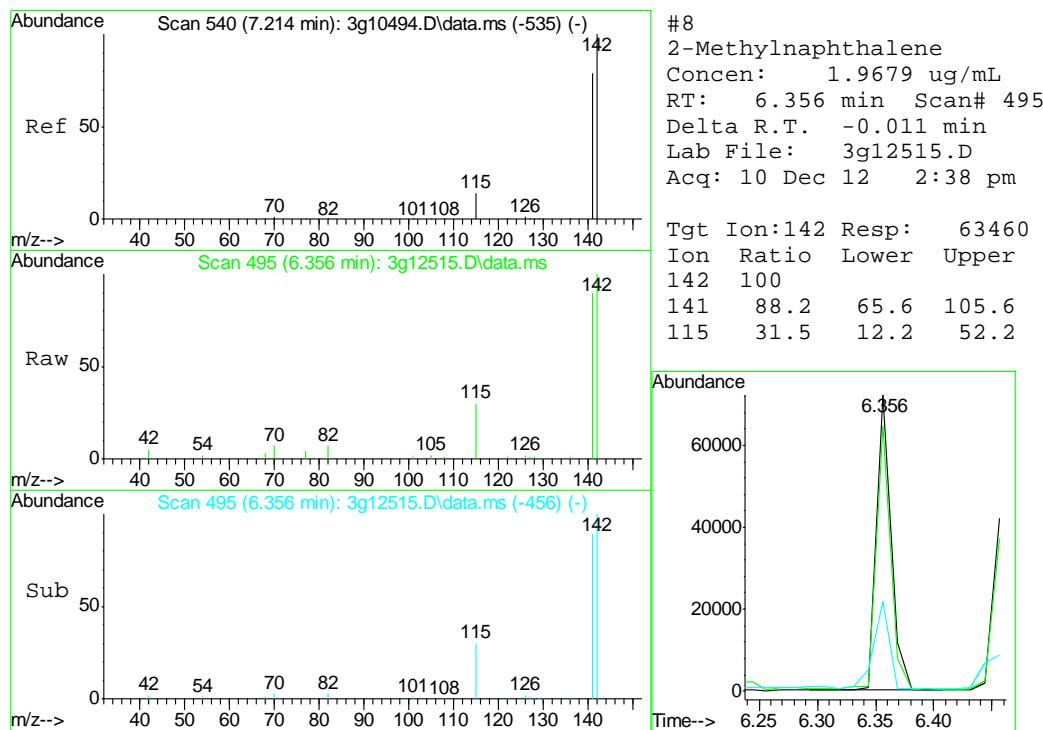
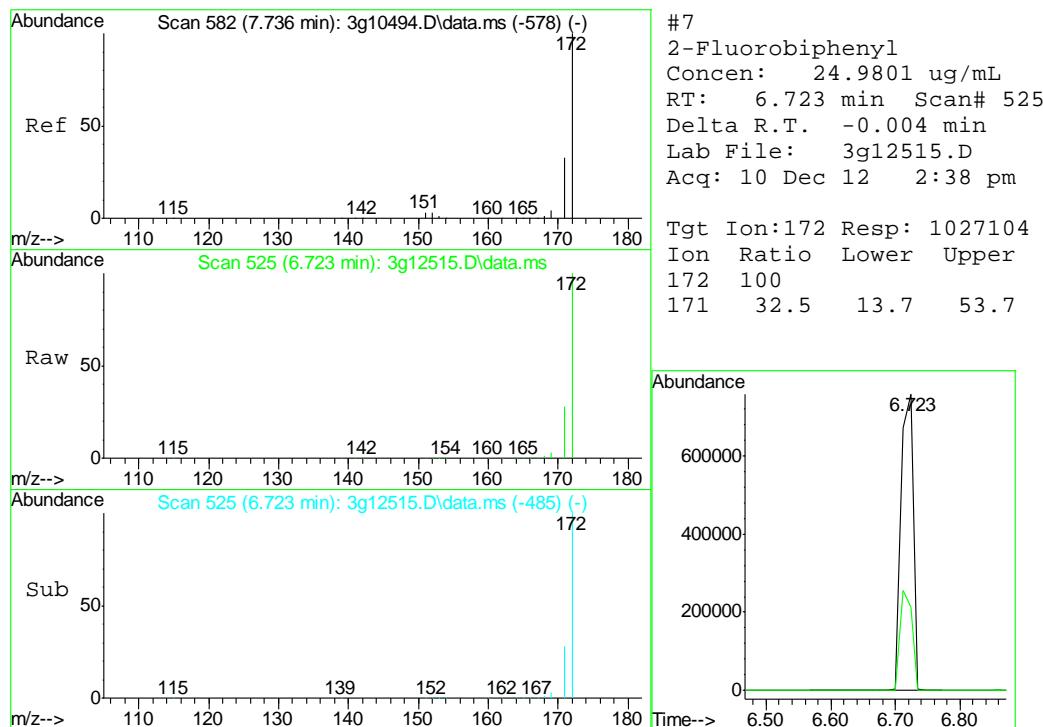
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 QLast Update : Tue Dec 04 08:50:28 2012
 Response via : Initial Calibration

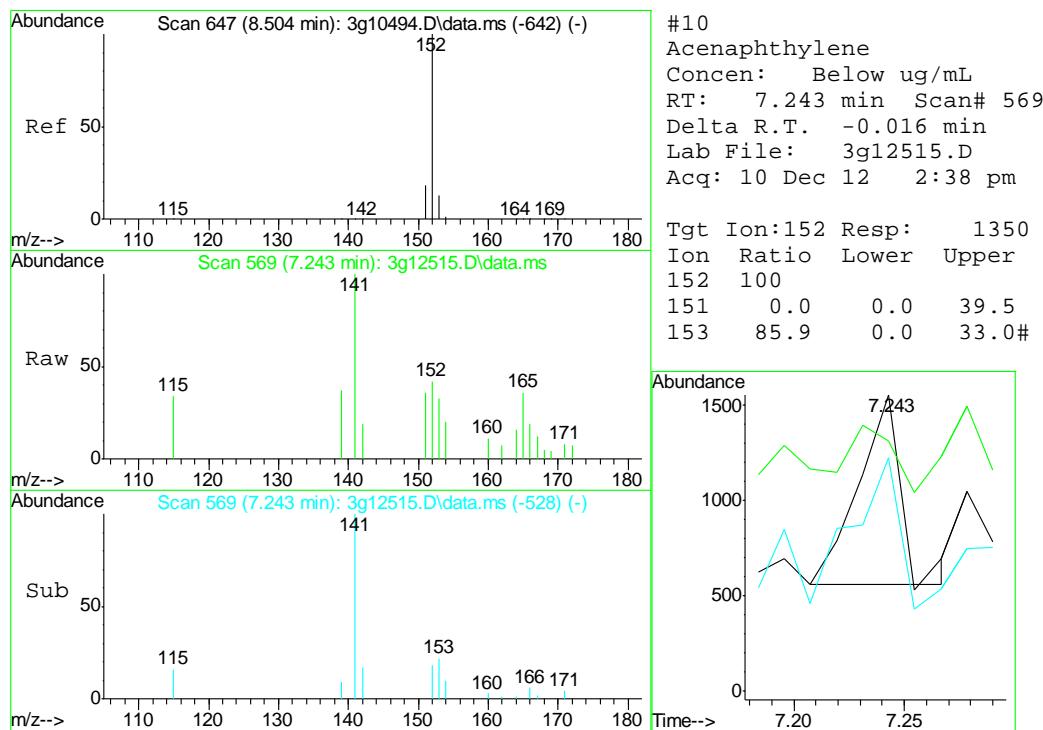
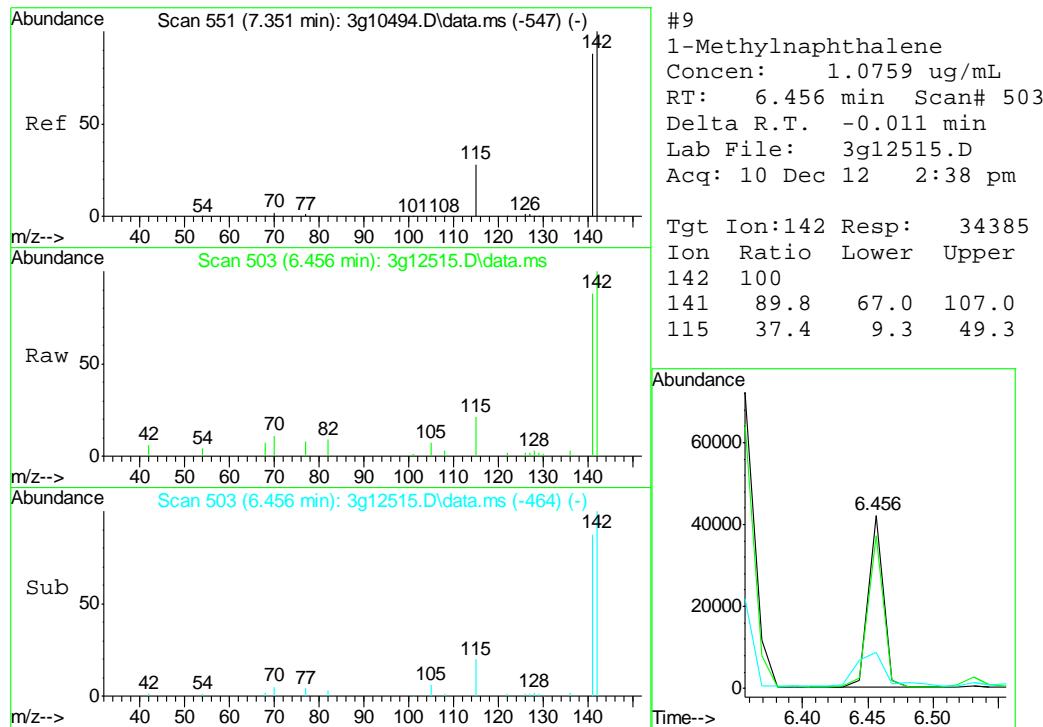


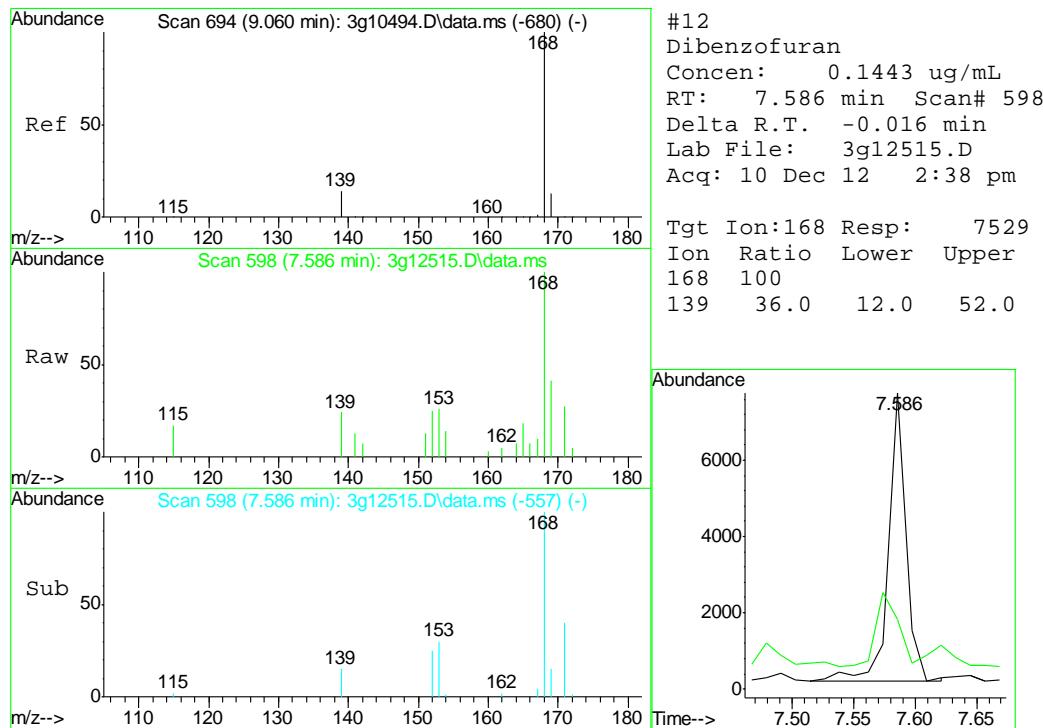
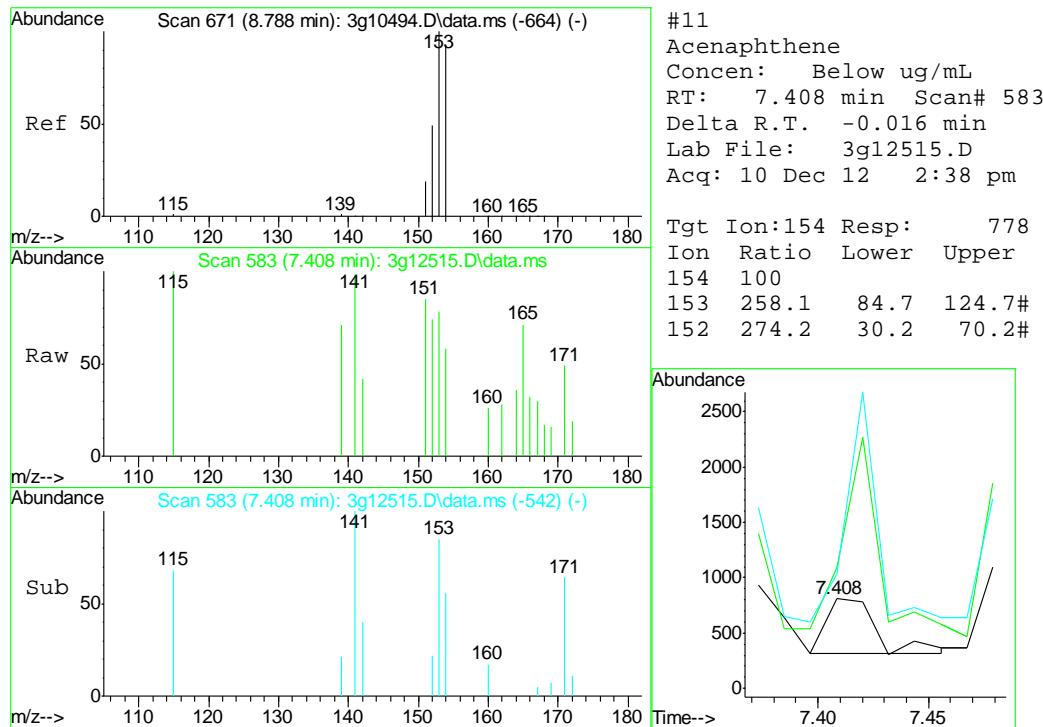


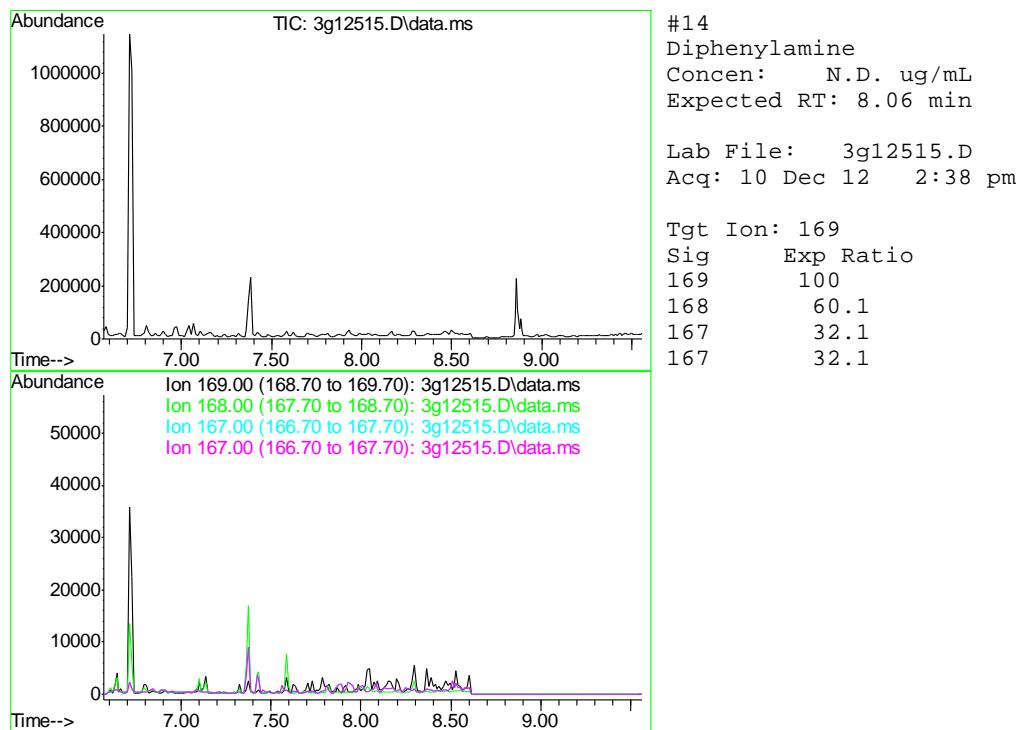
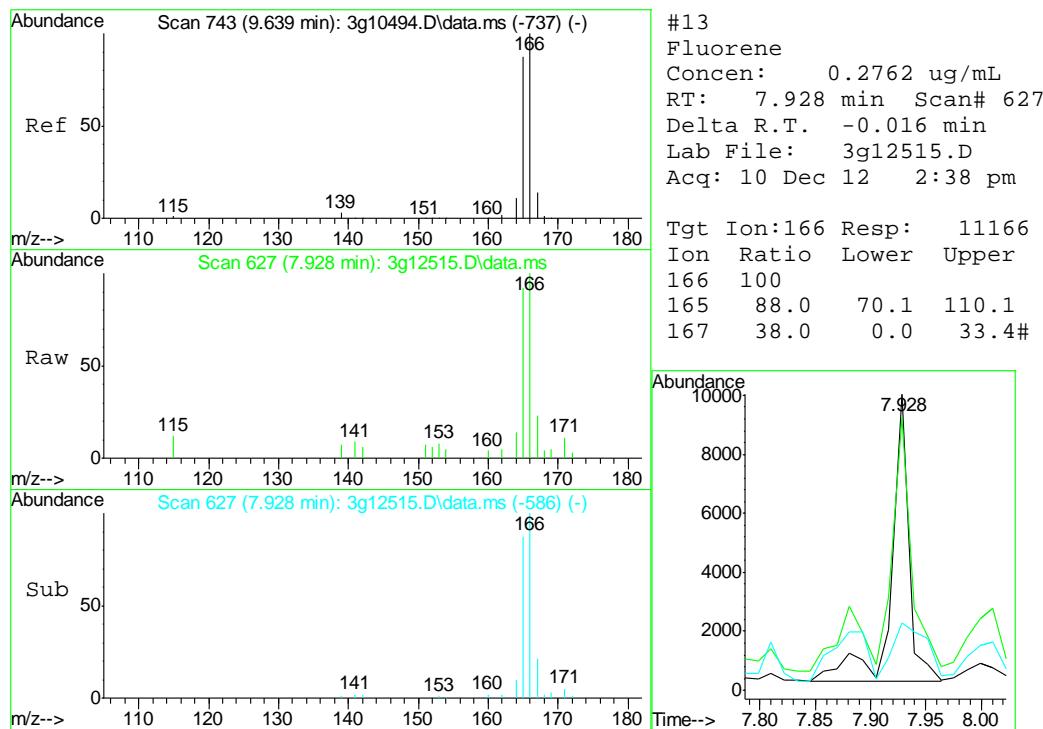


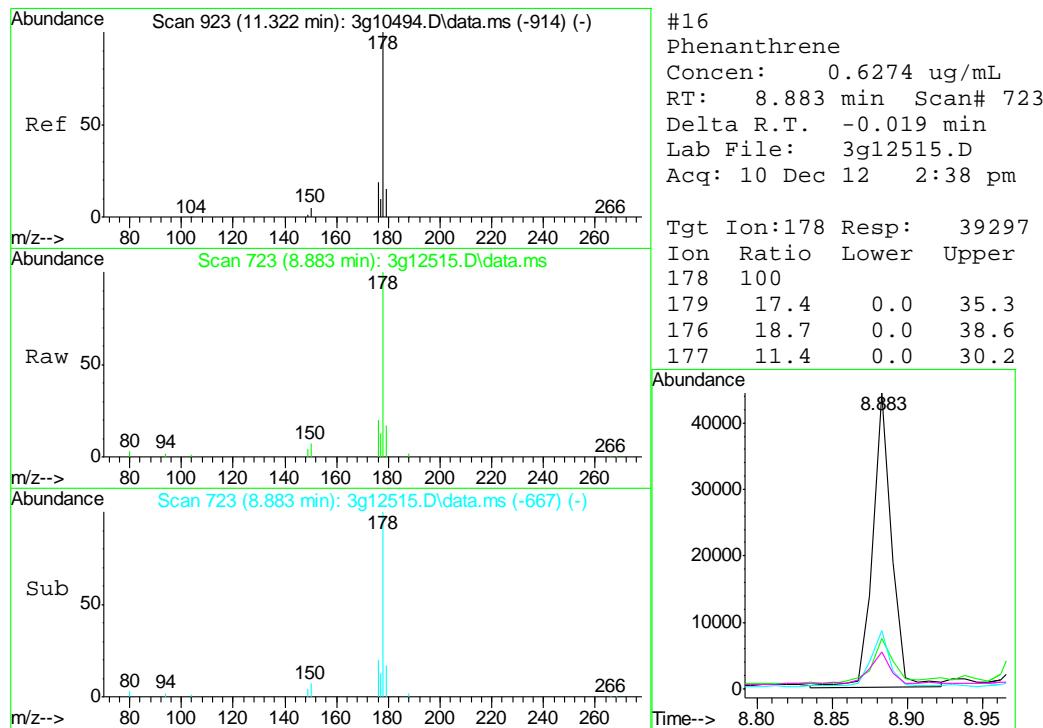
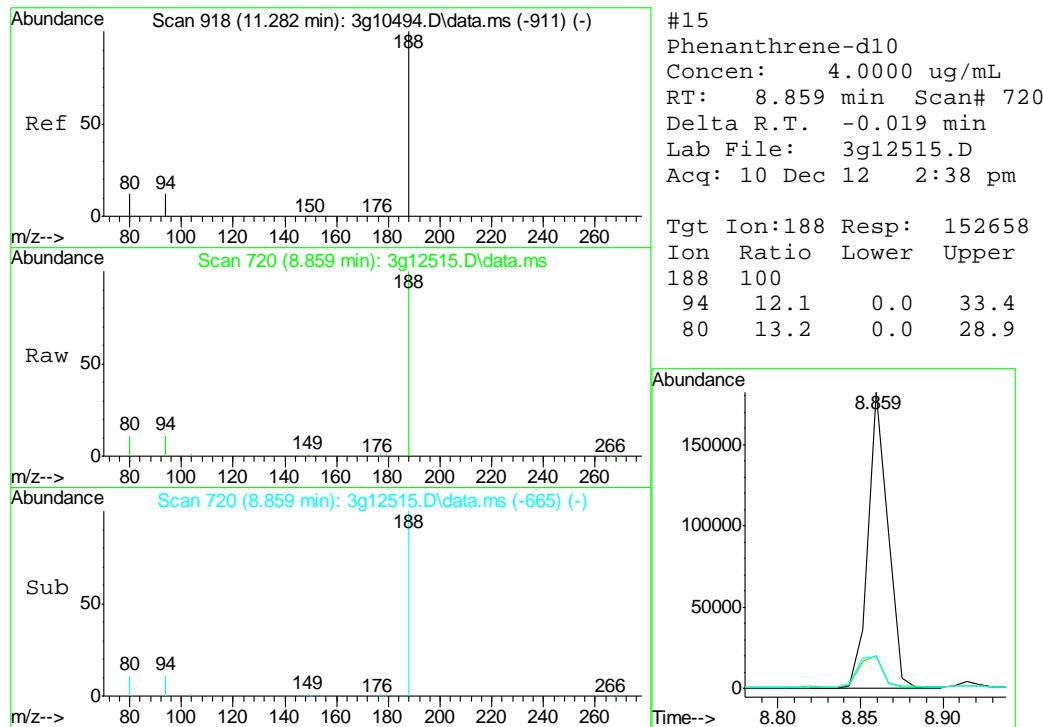


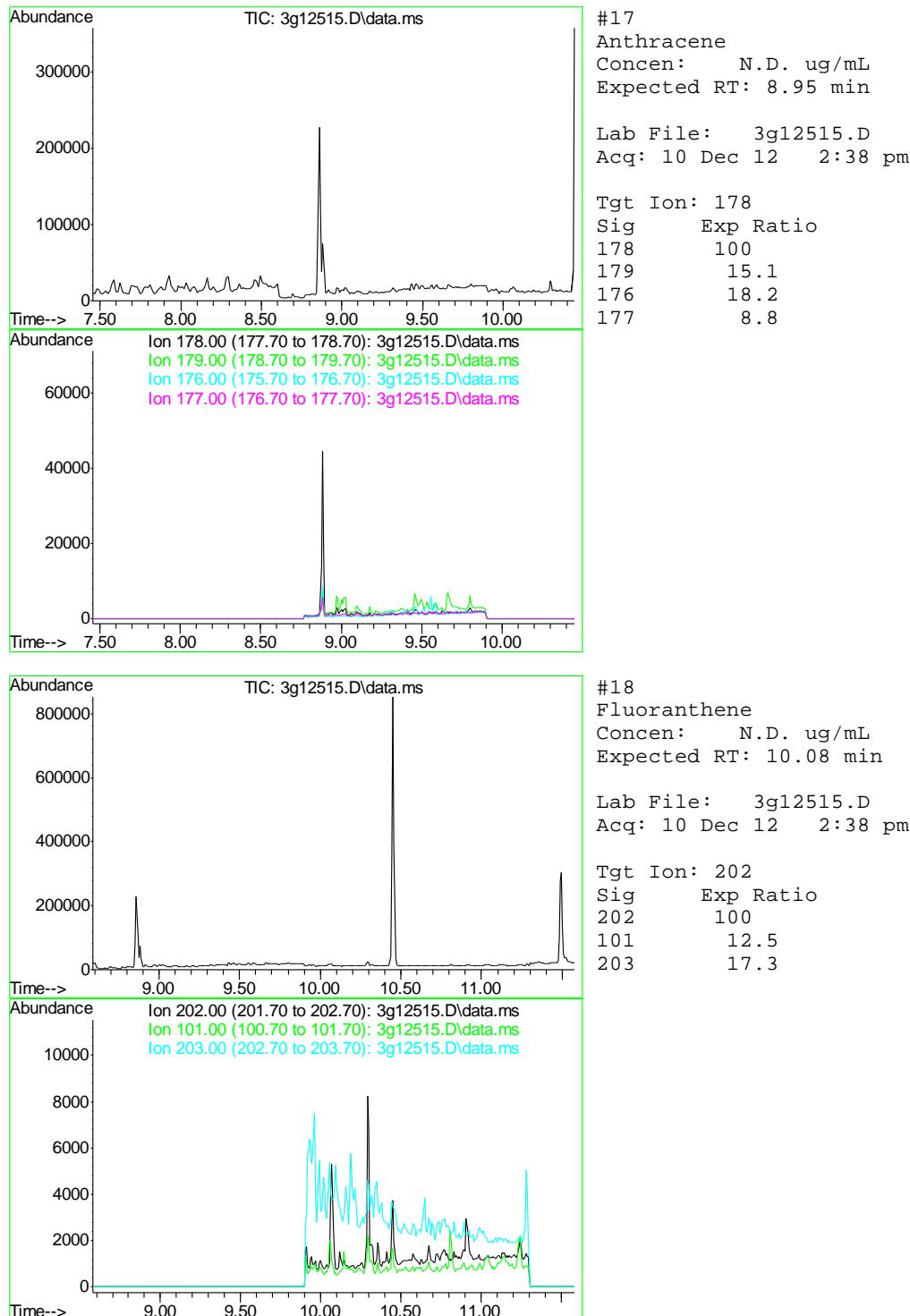


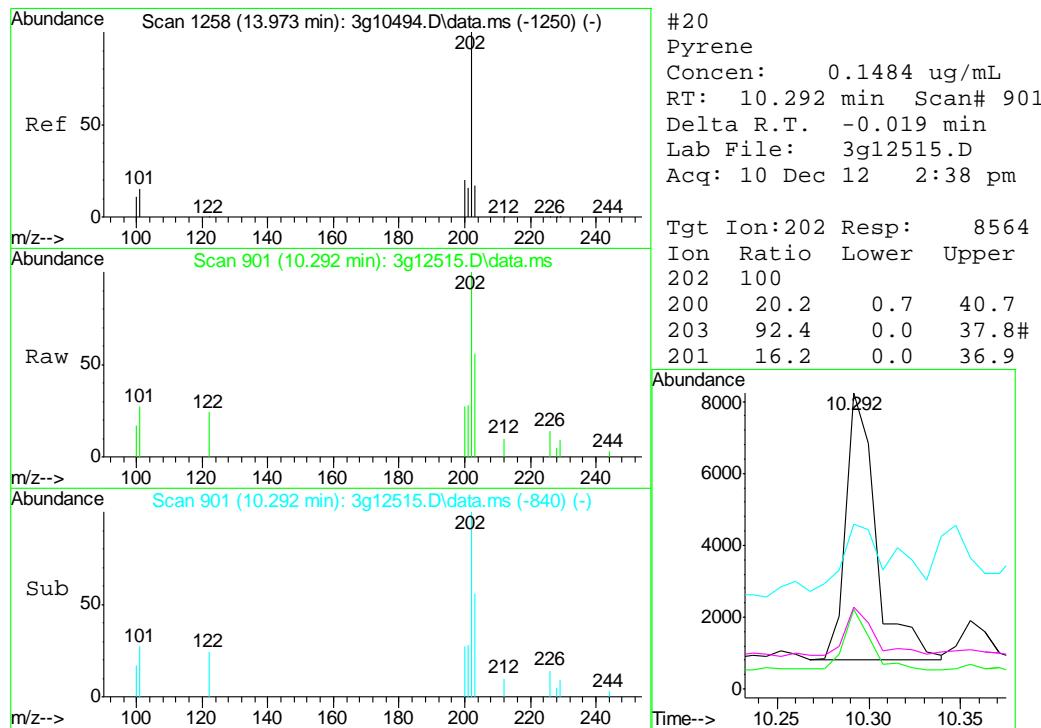
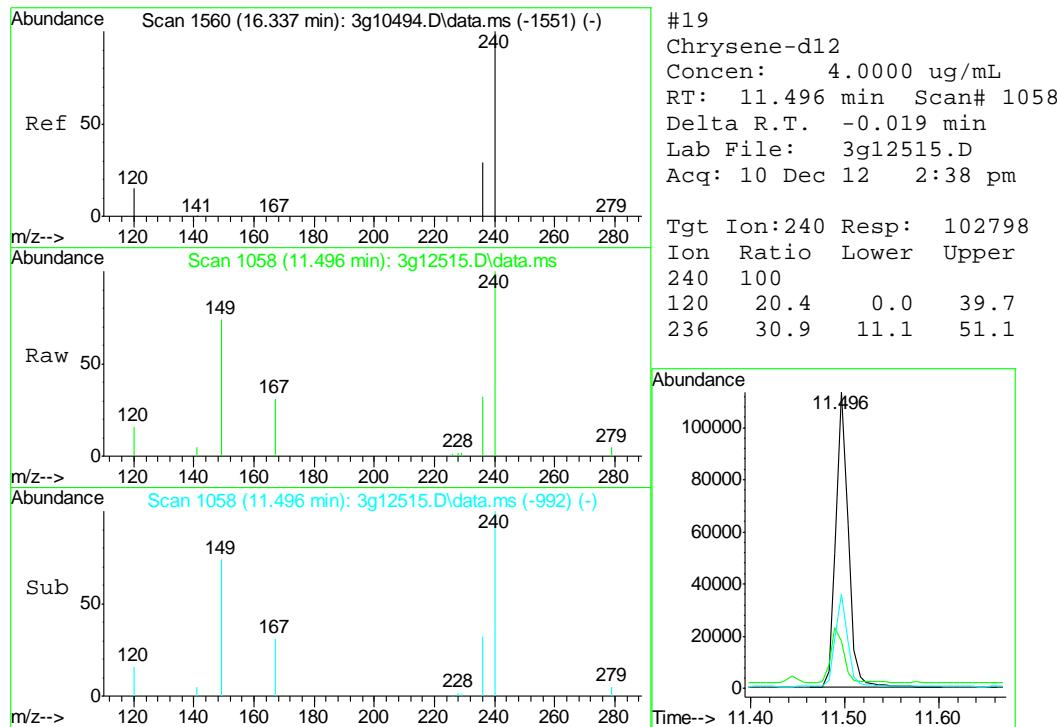


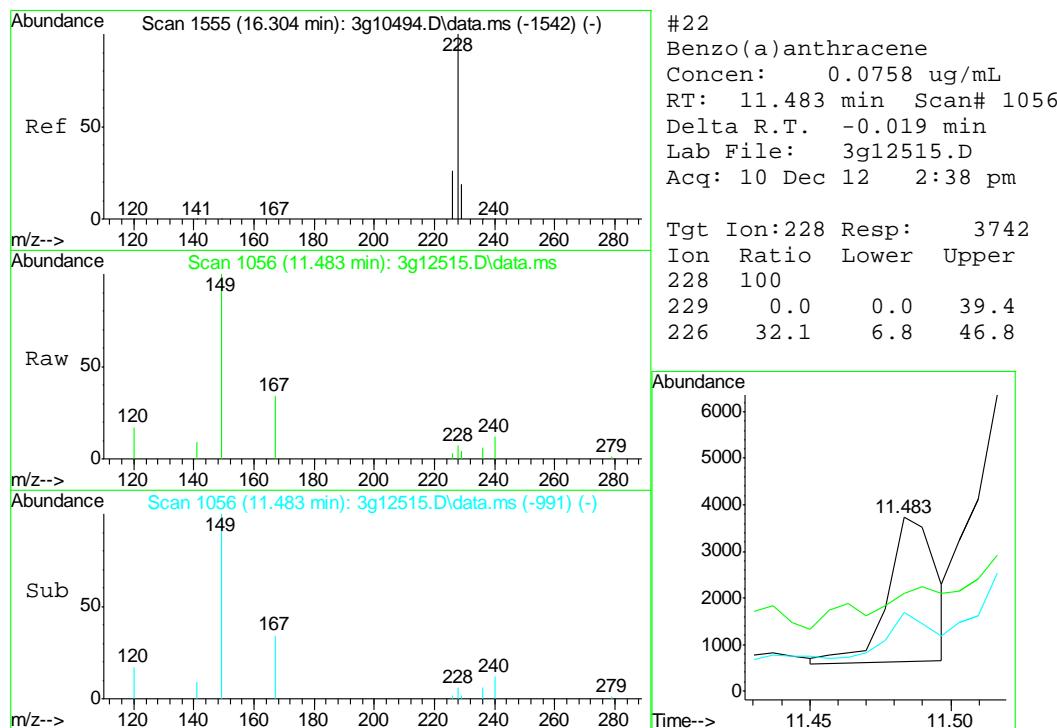
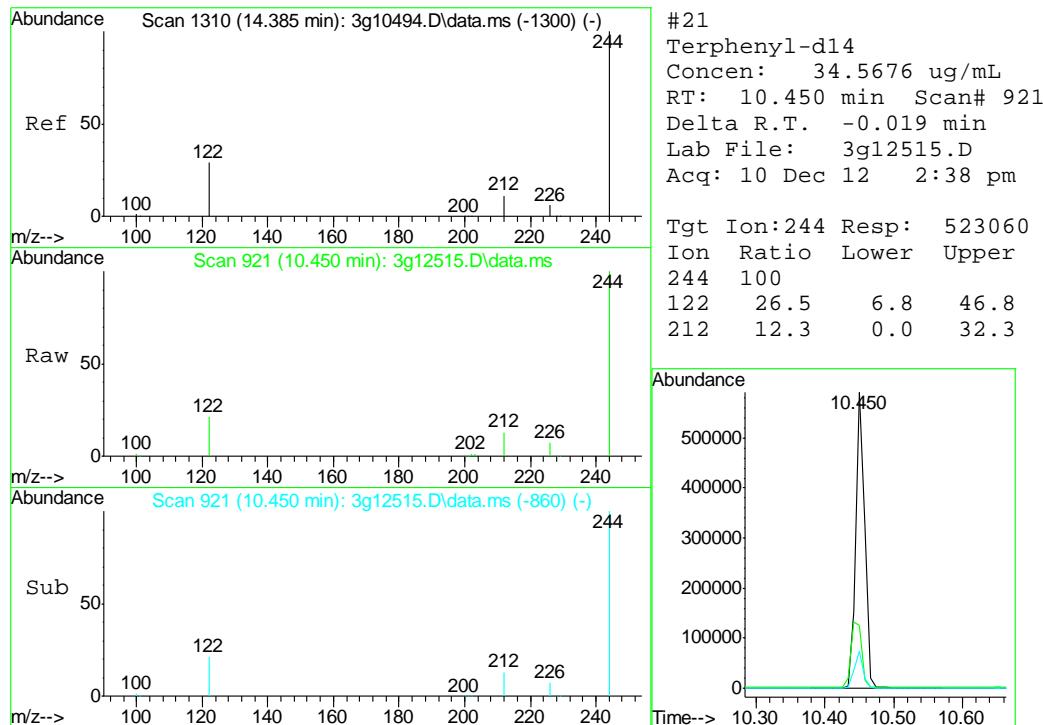


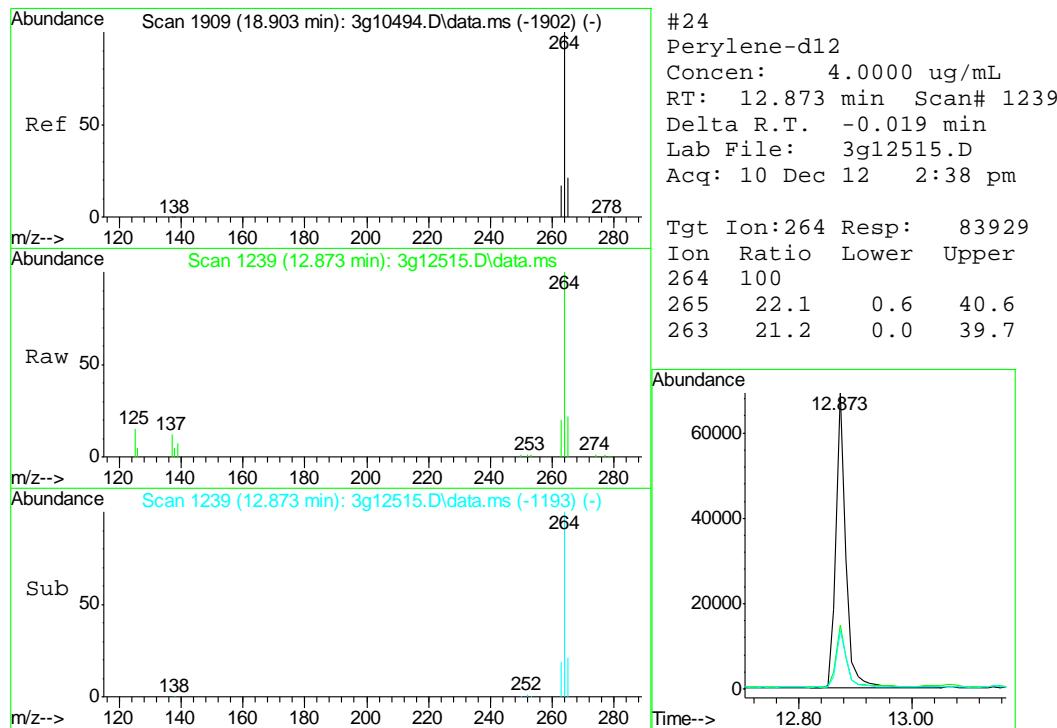
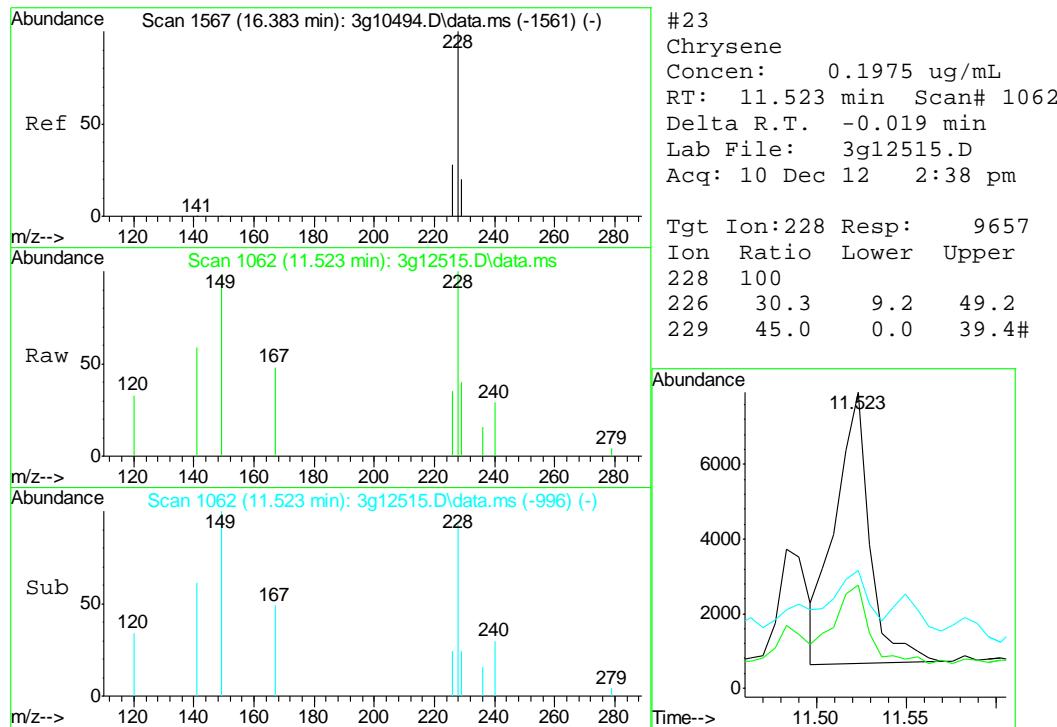


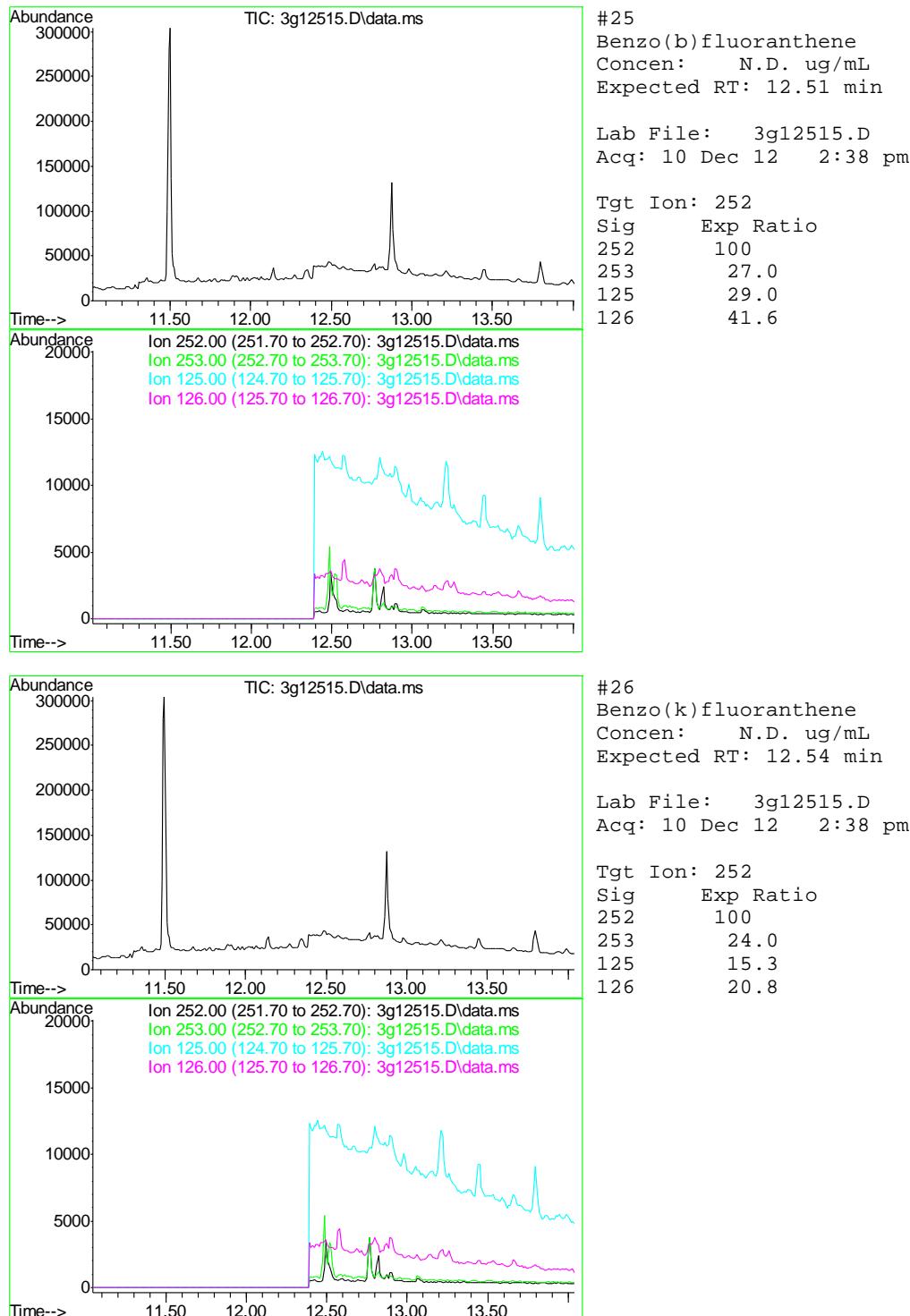


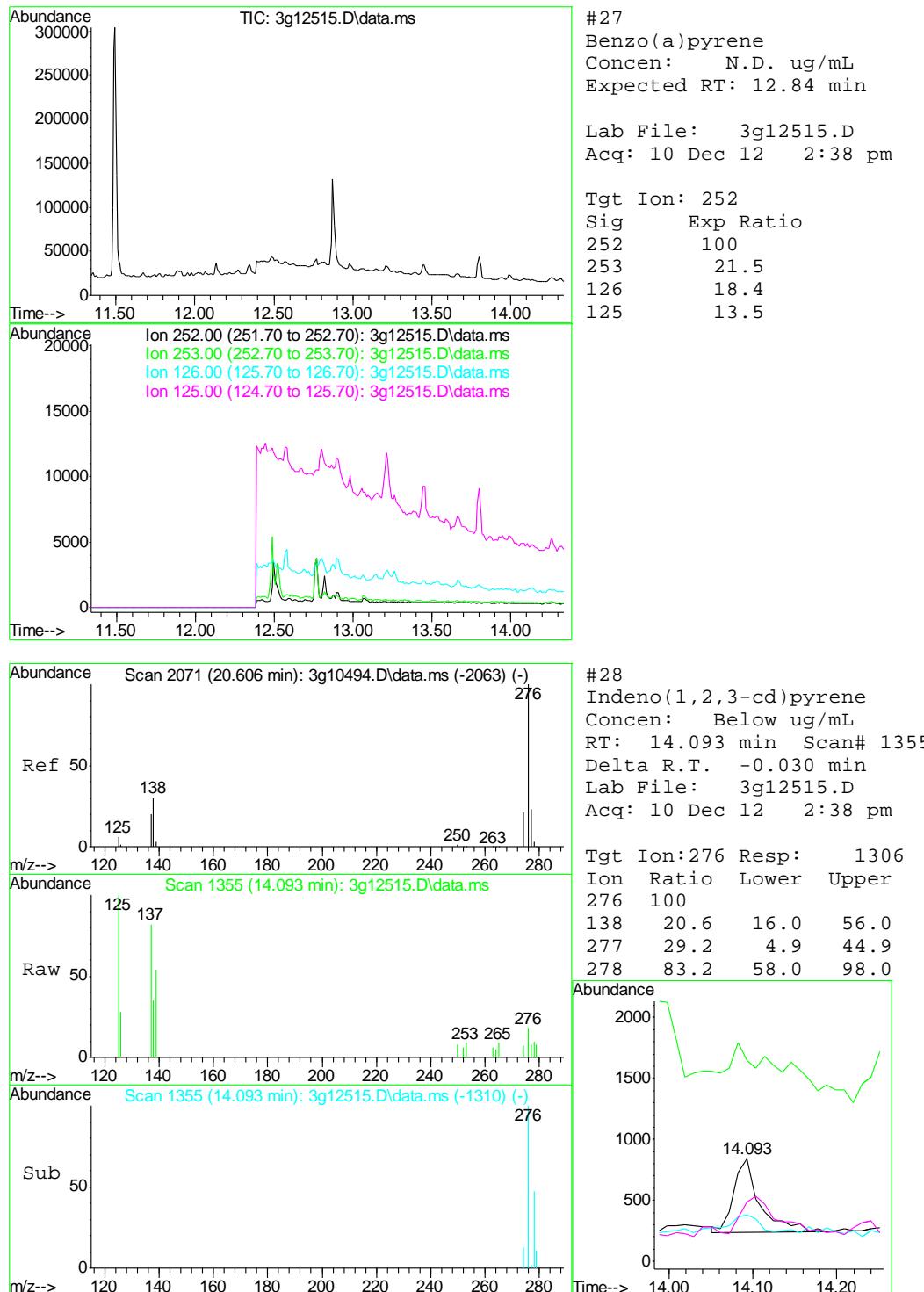


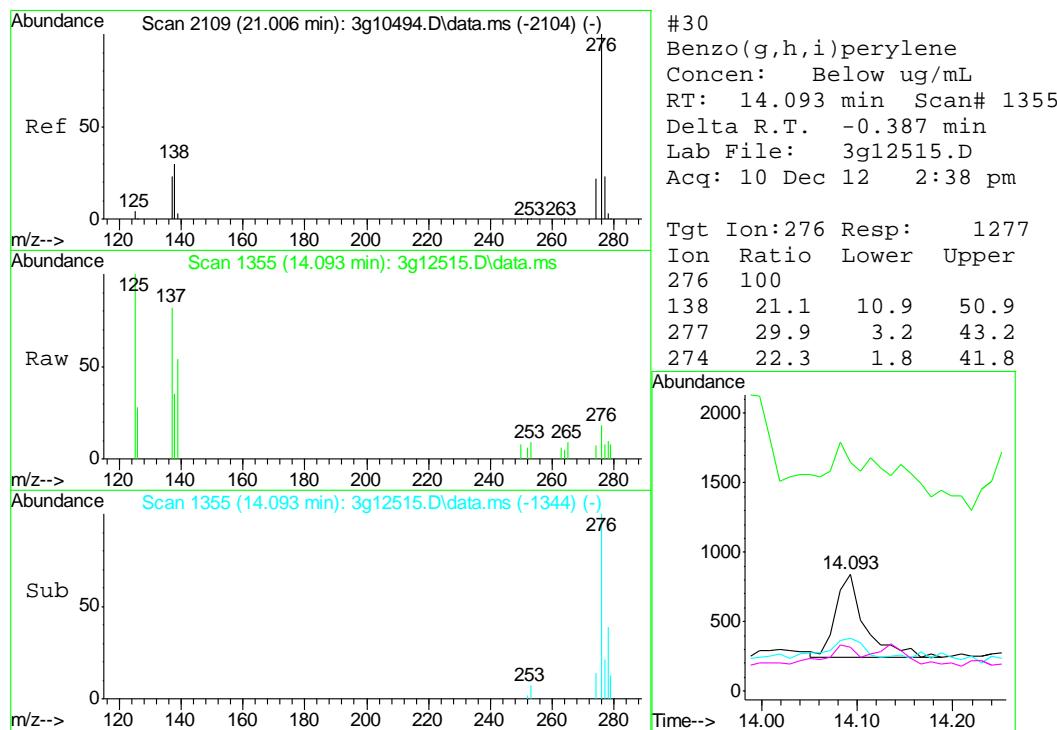
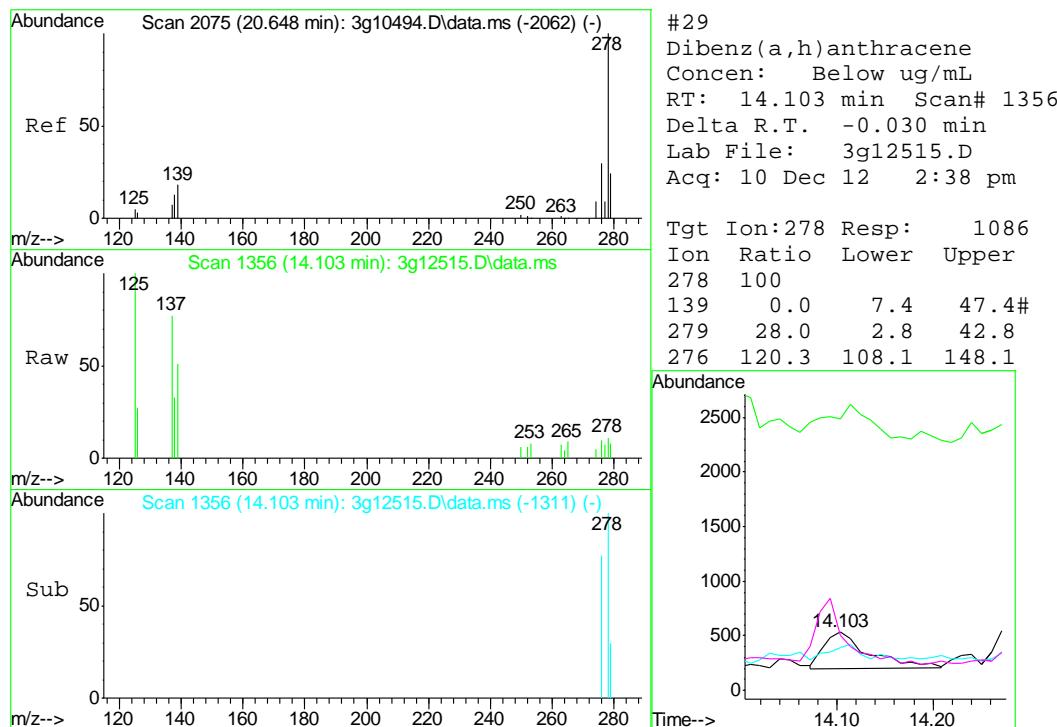












Quantitation Report (QT Reviewed)

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 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

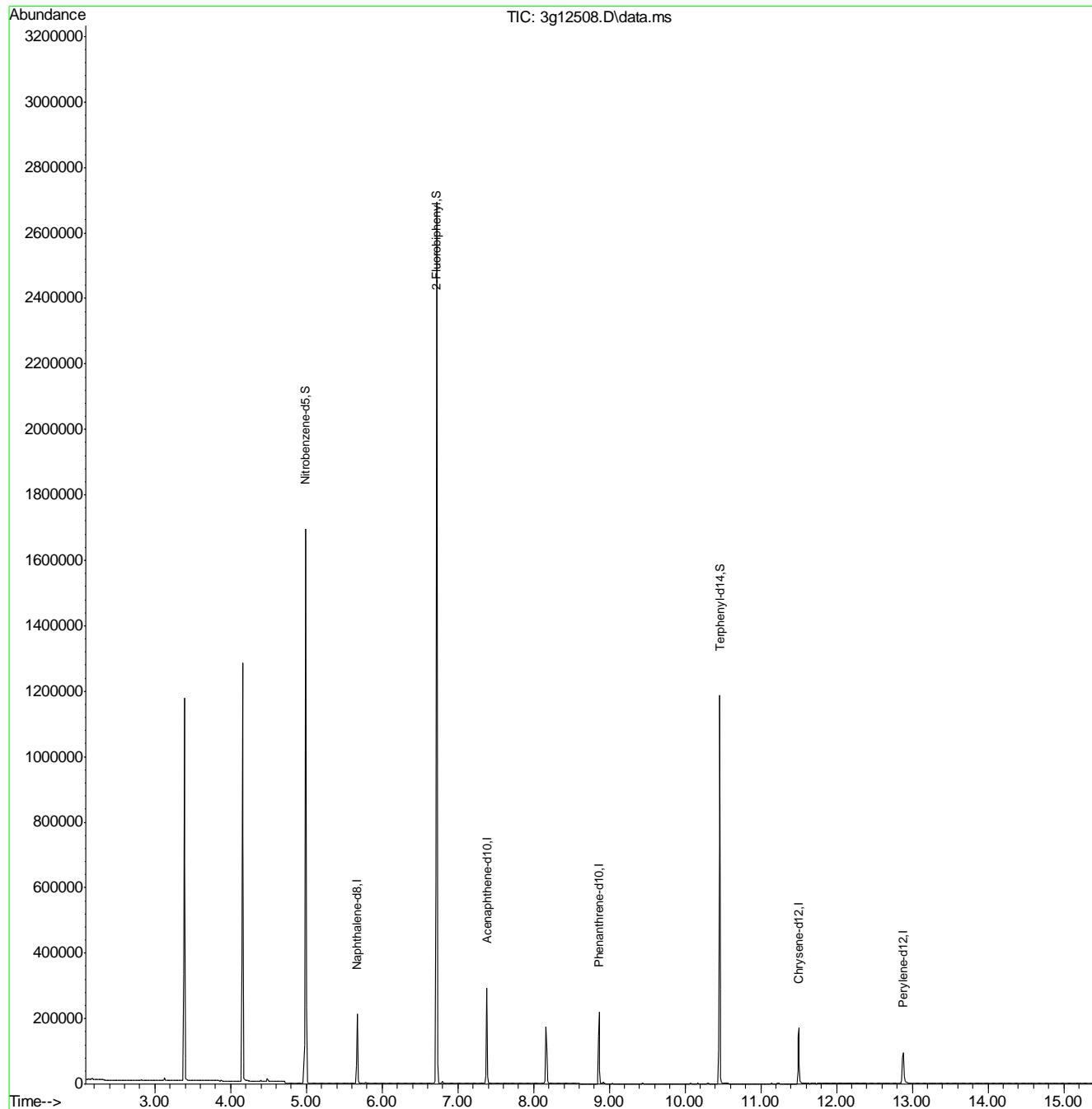
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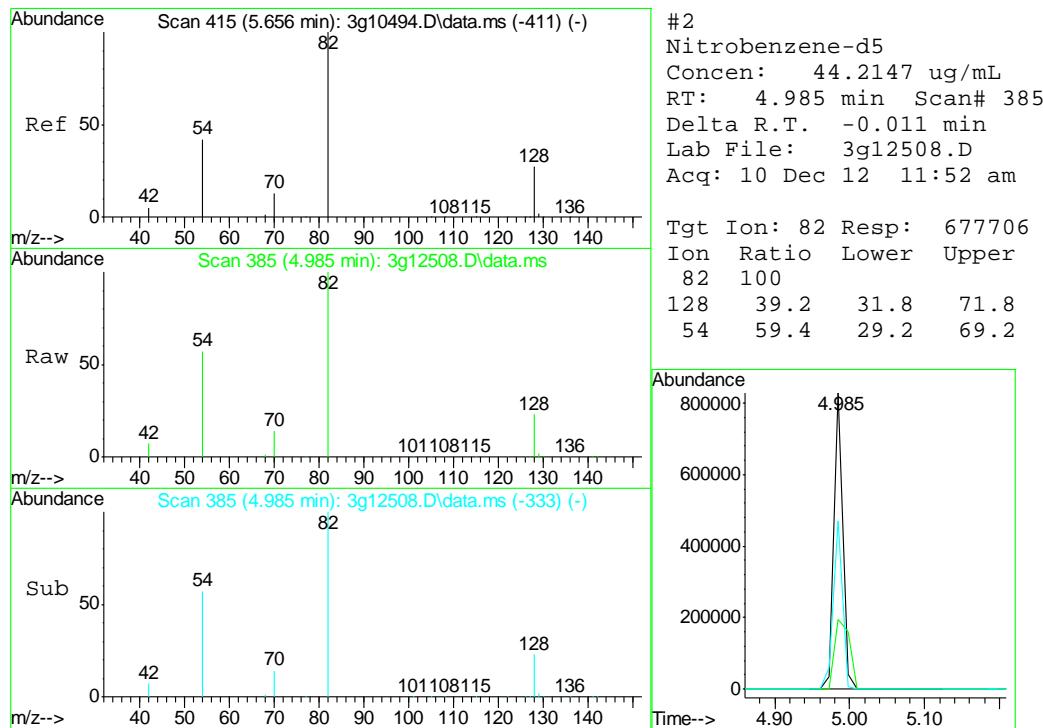
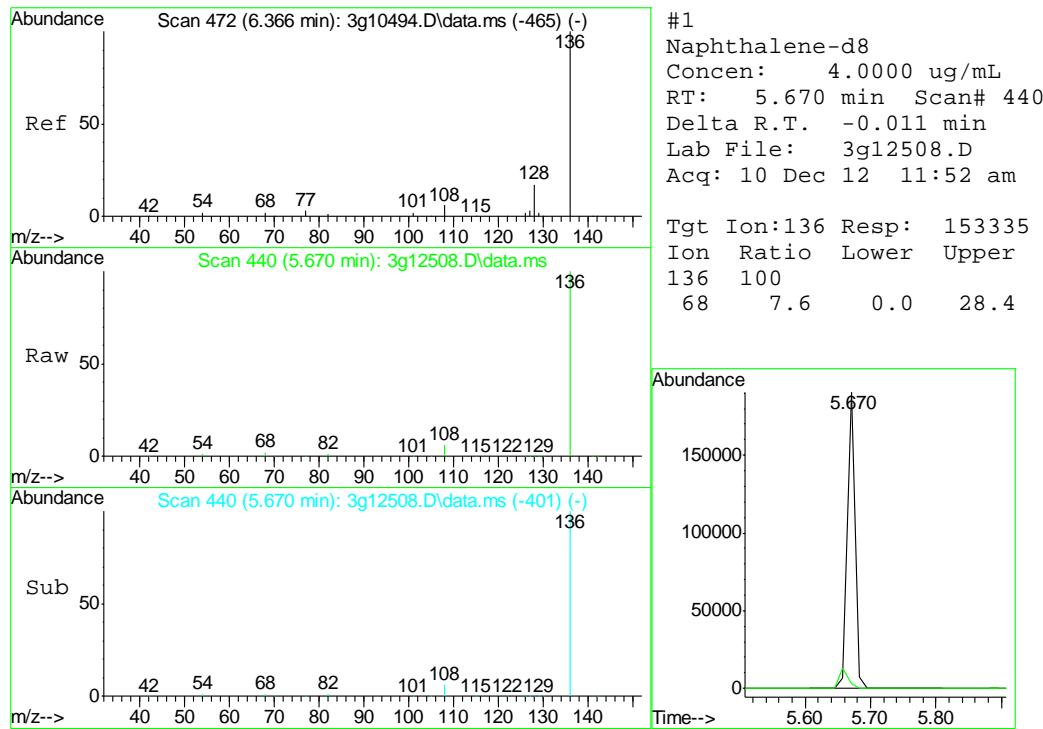
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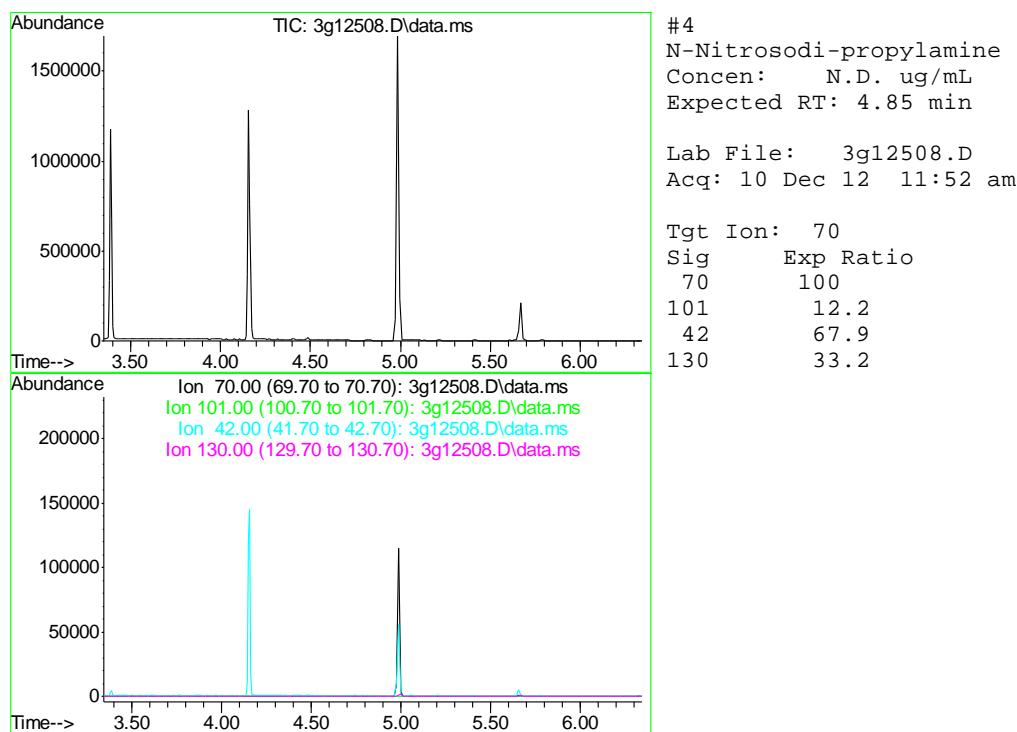
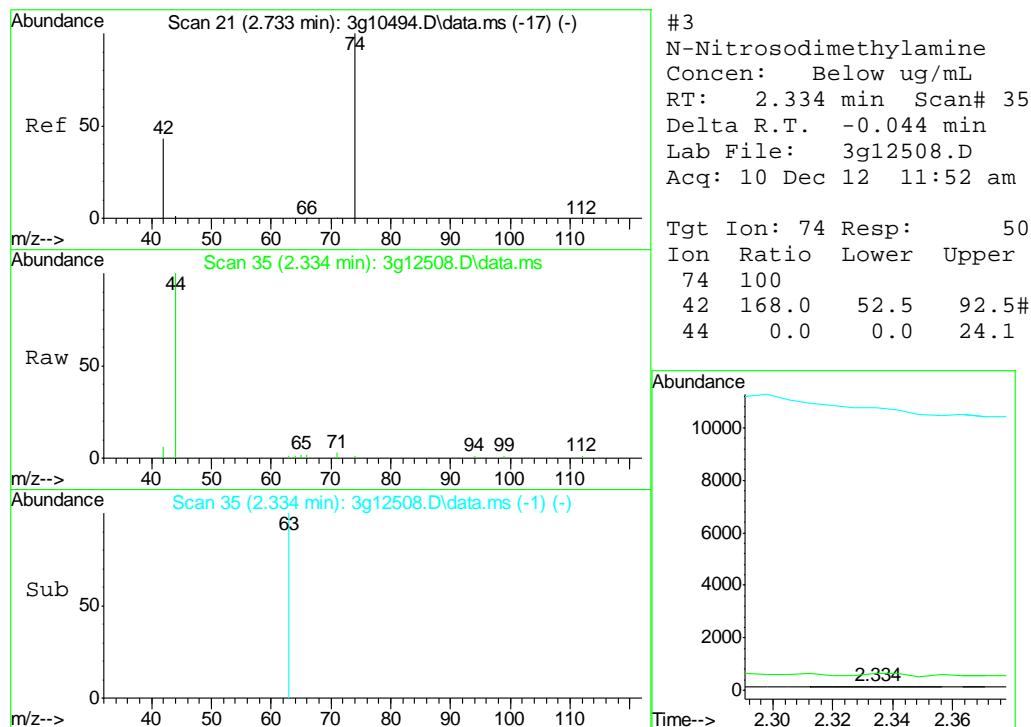
Quantitation Report (QT Reviewed)

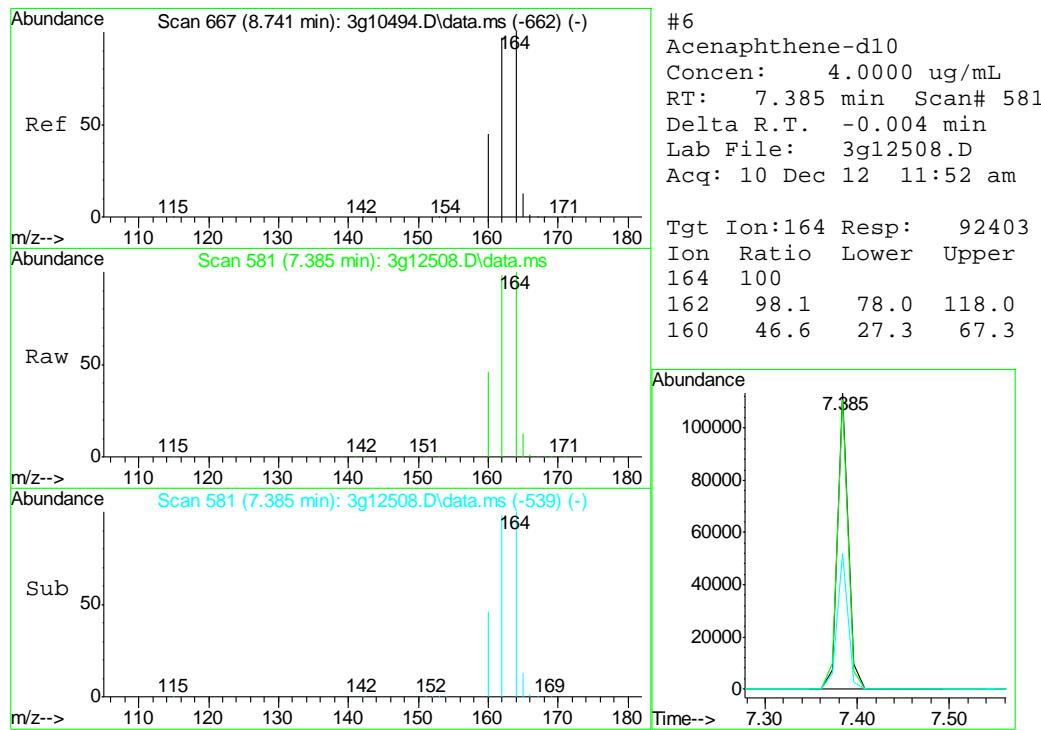
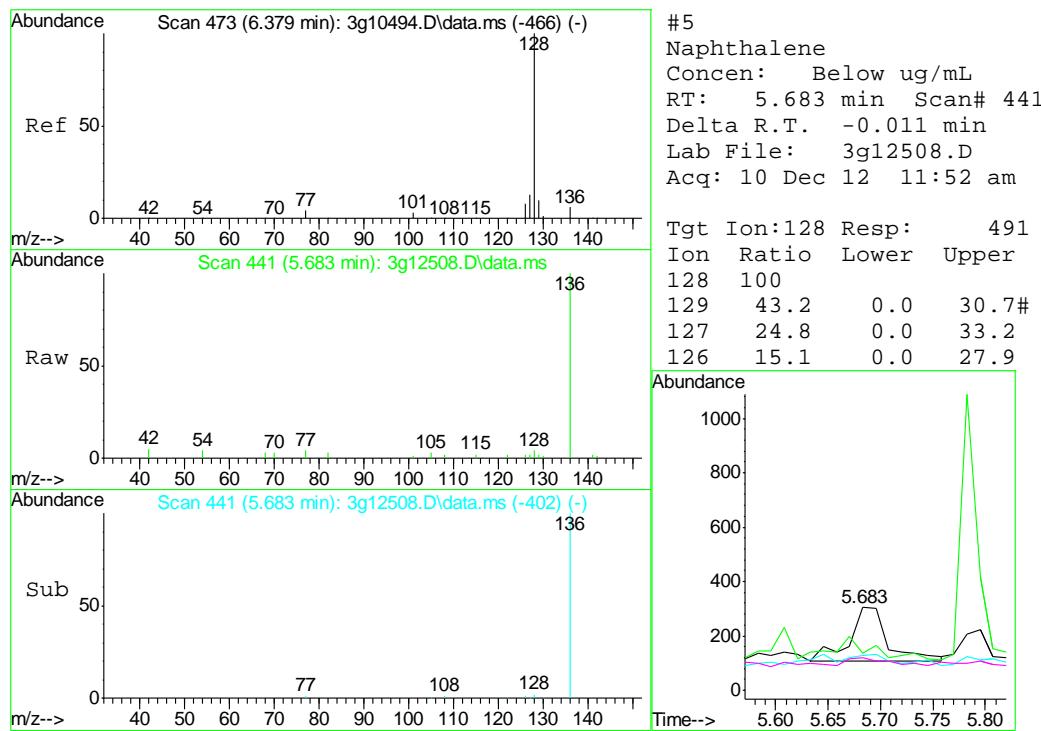
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 ALS Vial : 4 Sample Multiplier: 1

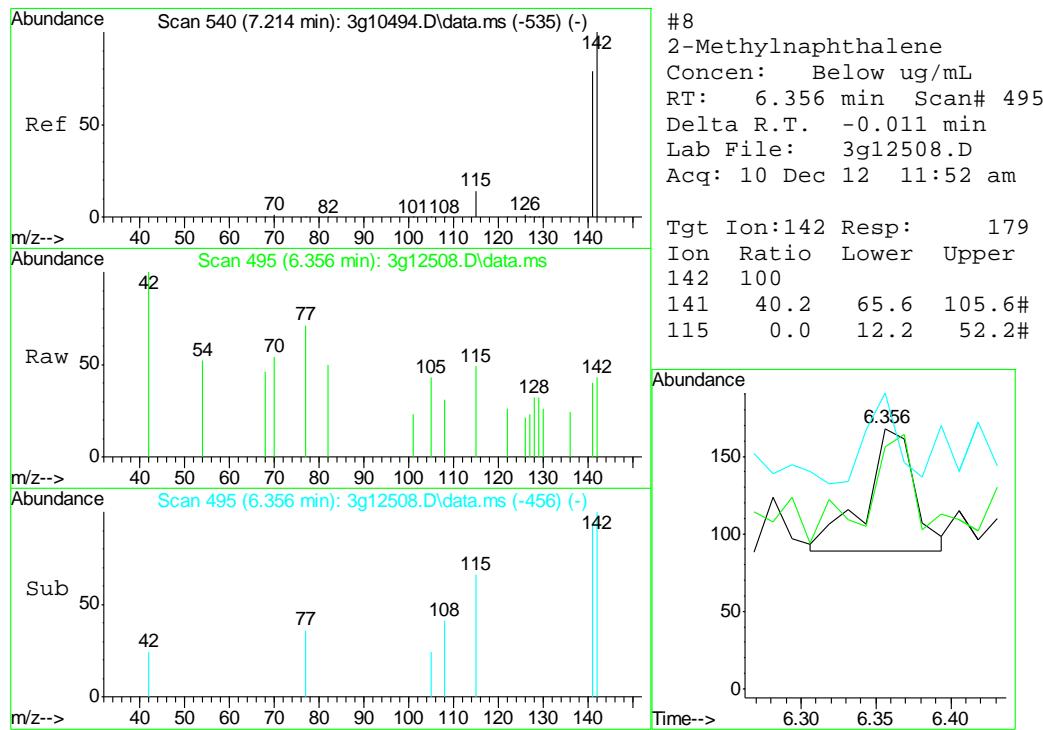
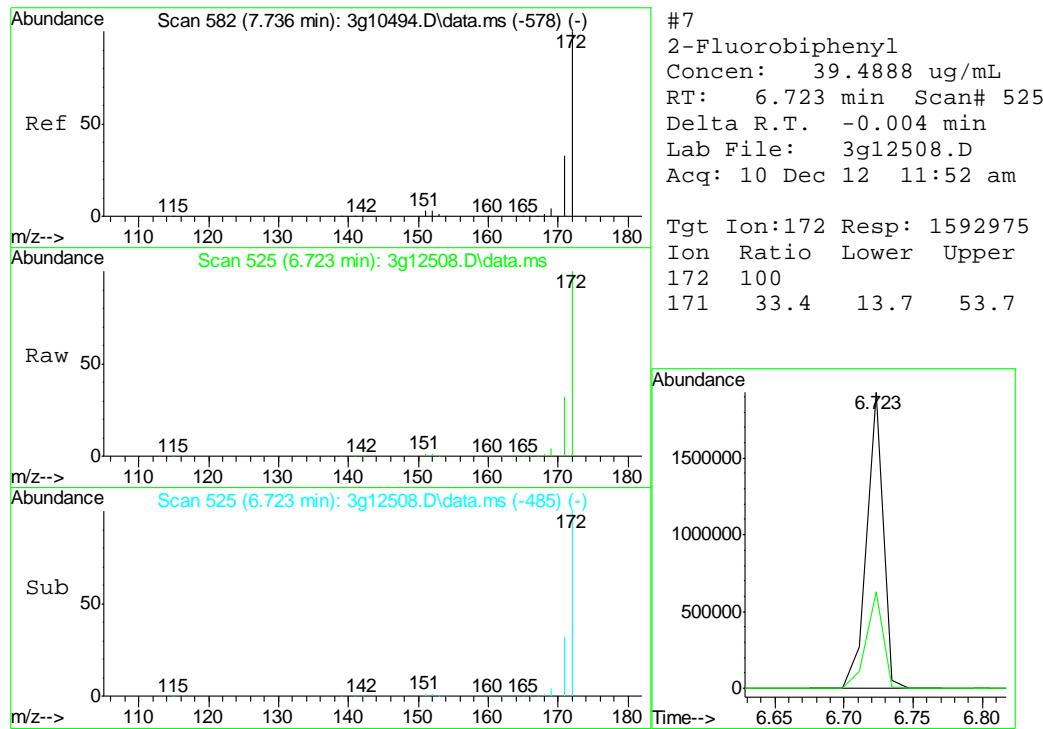
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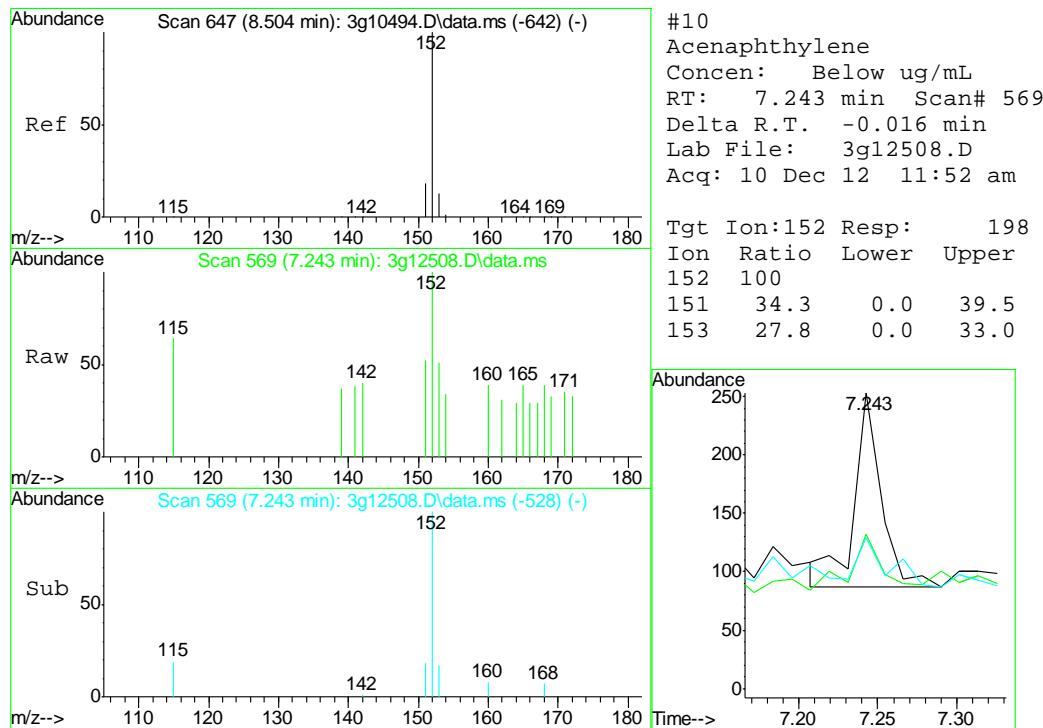
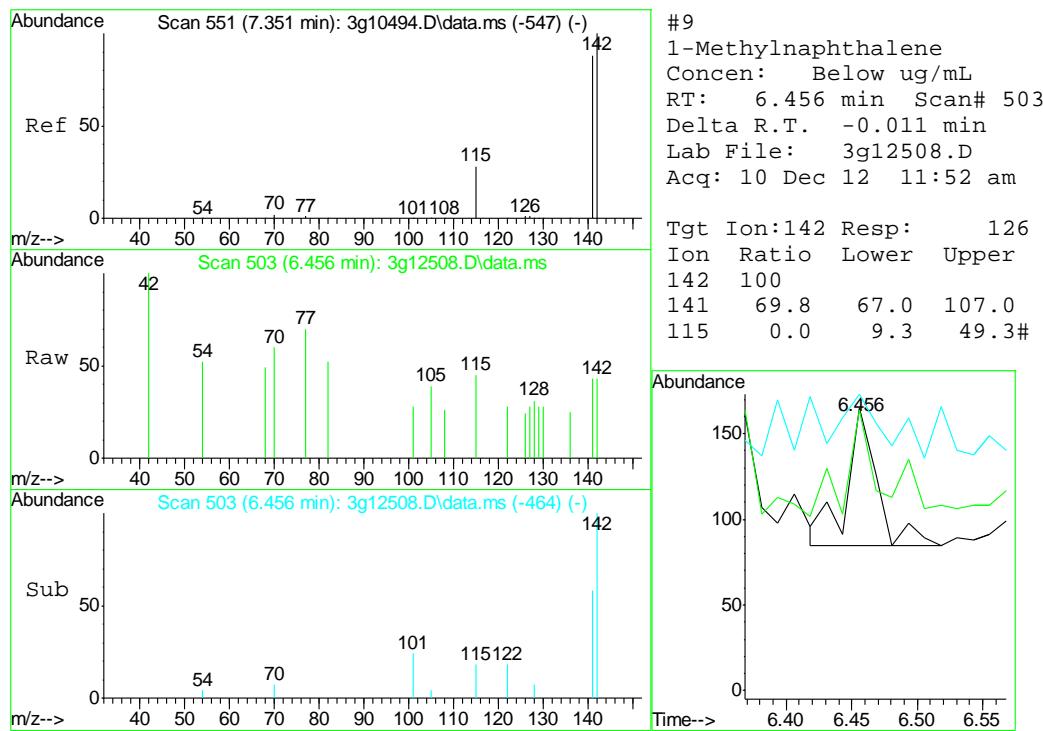


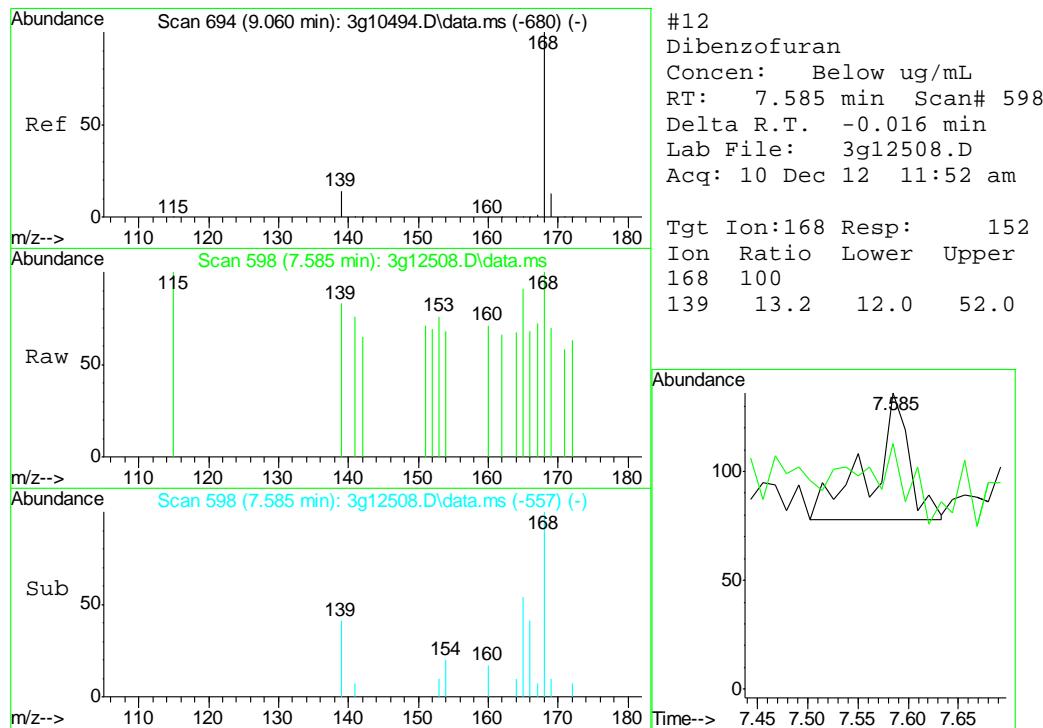
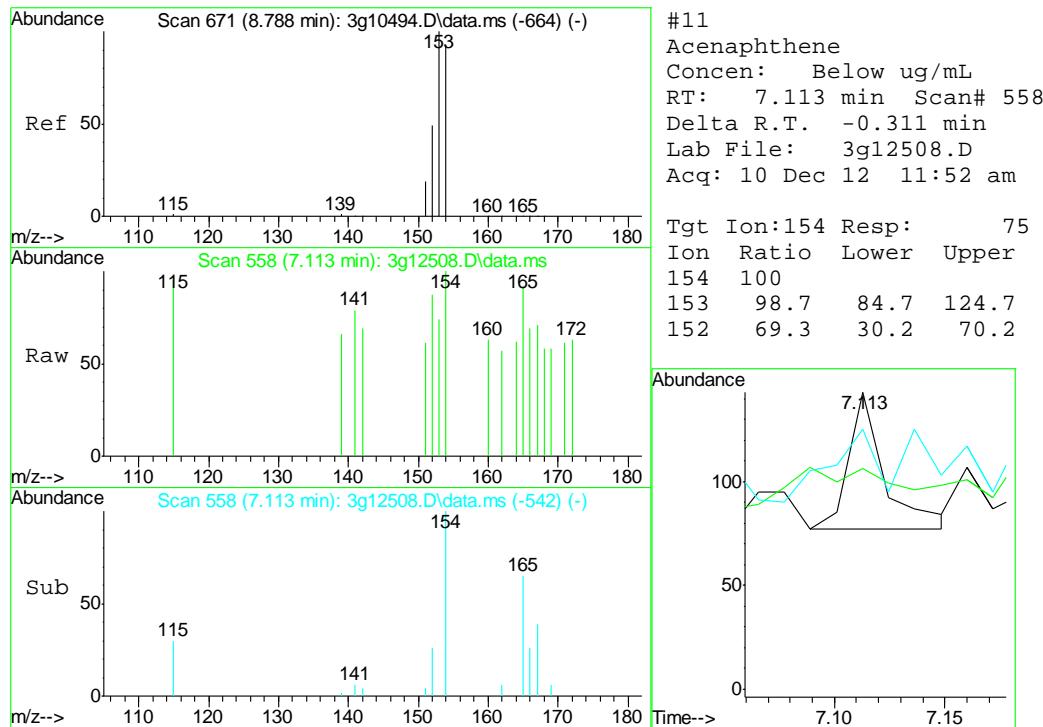


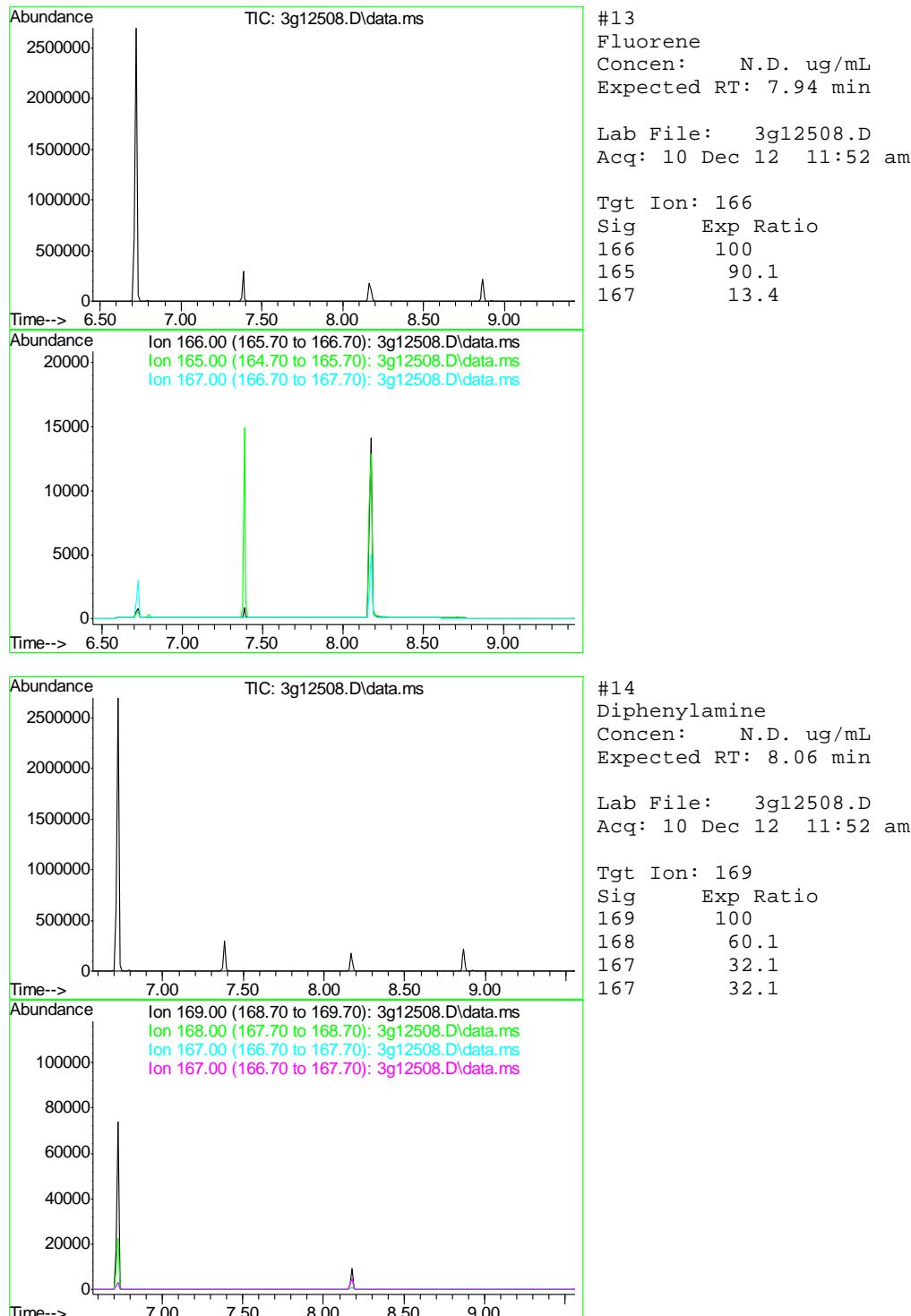


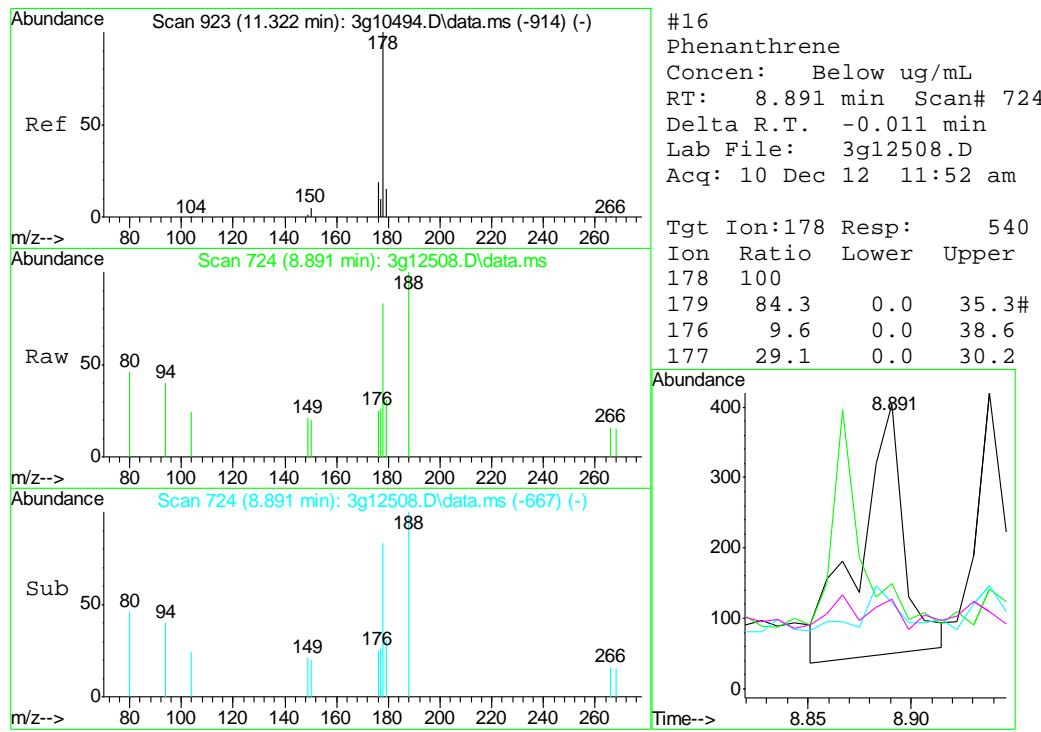
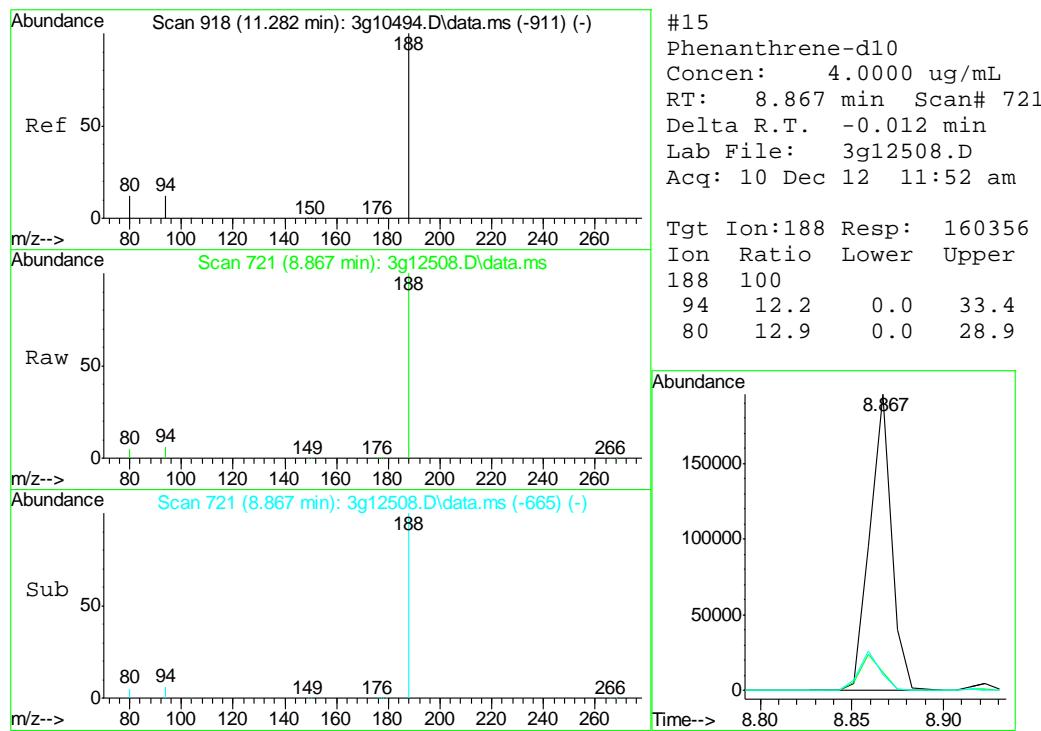


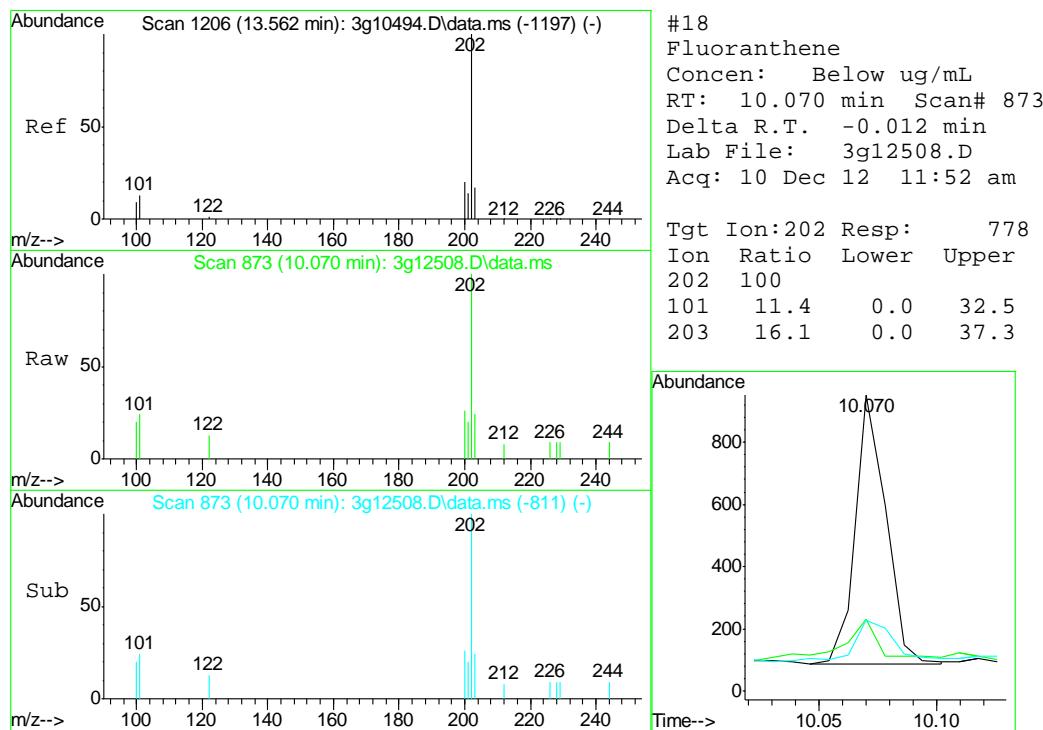
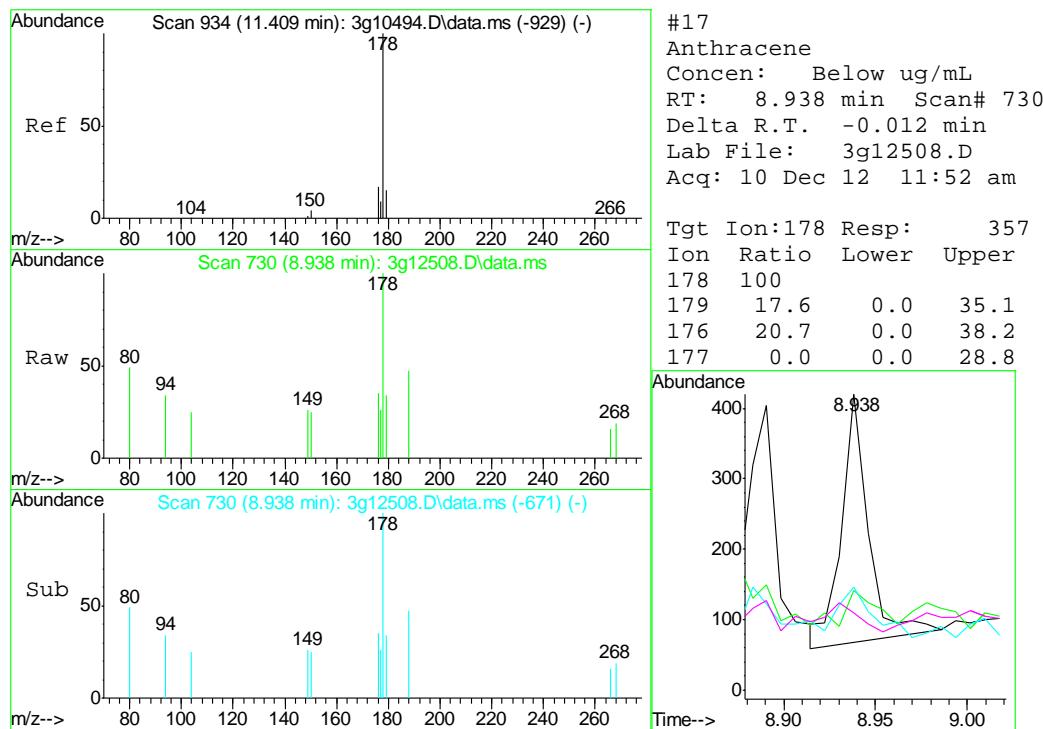


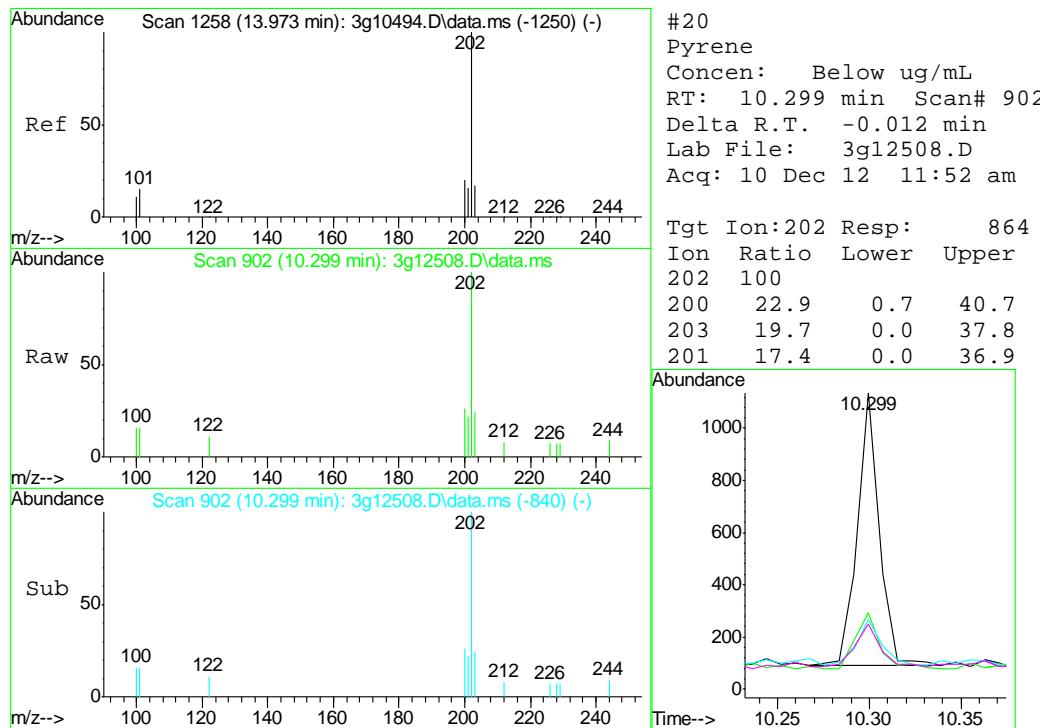
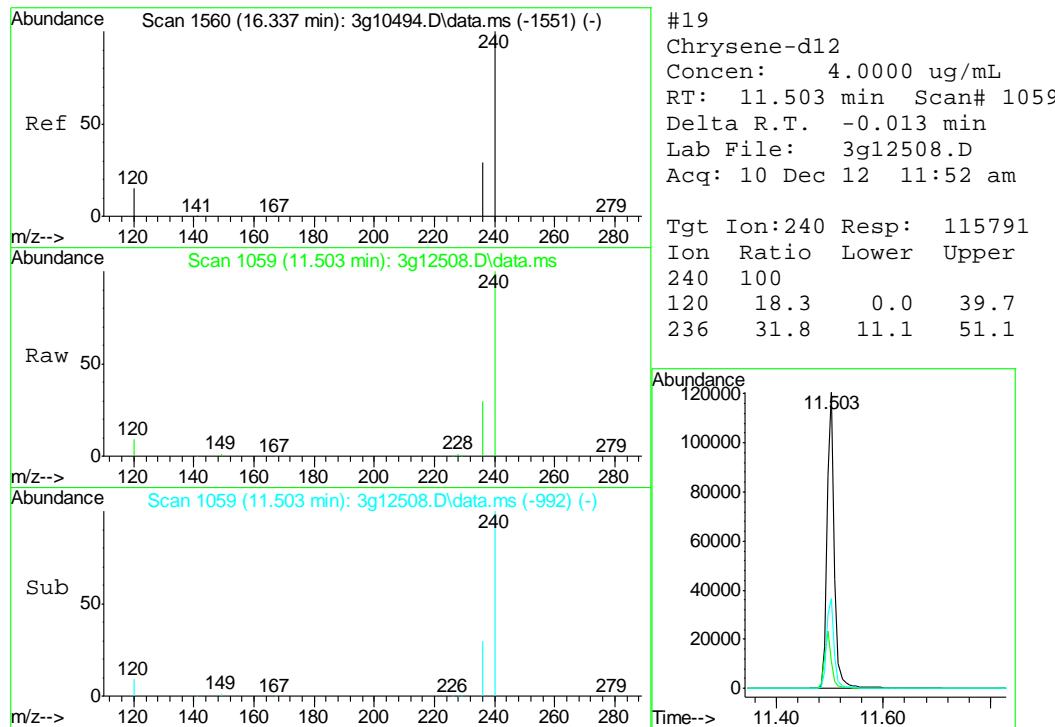


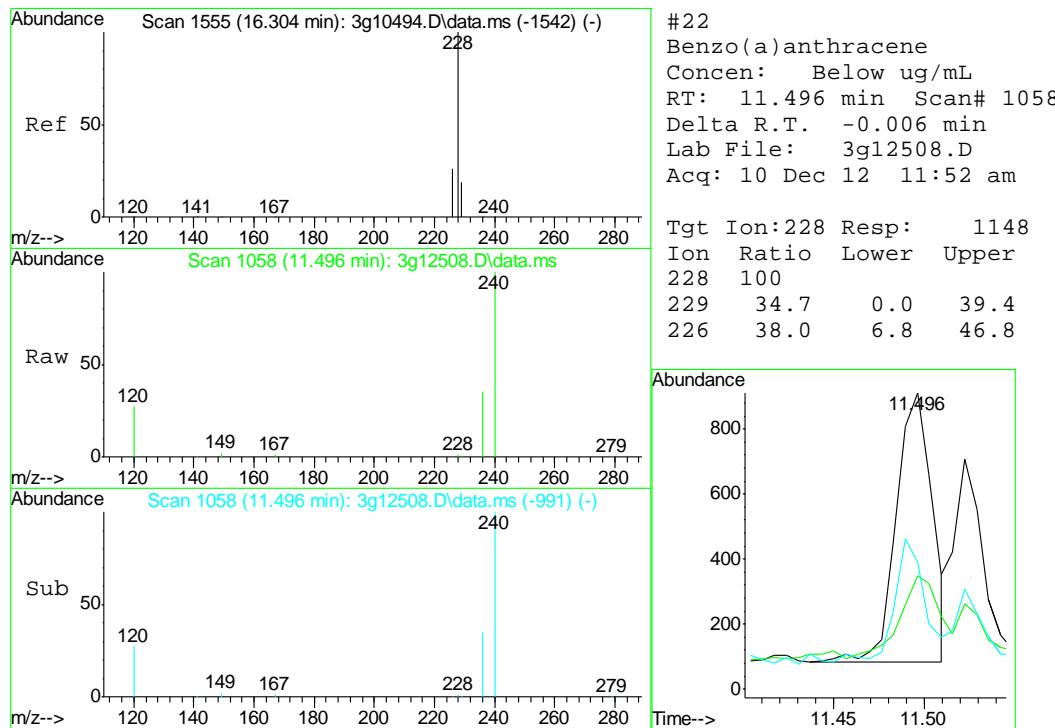
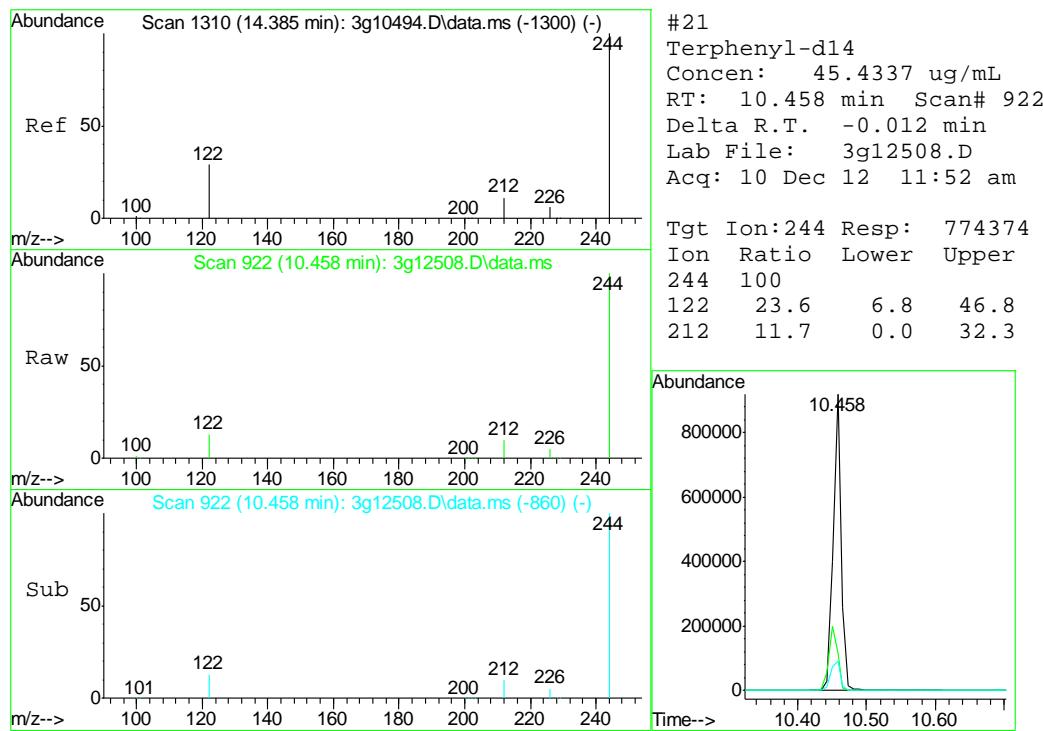


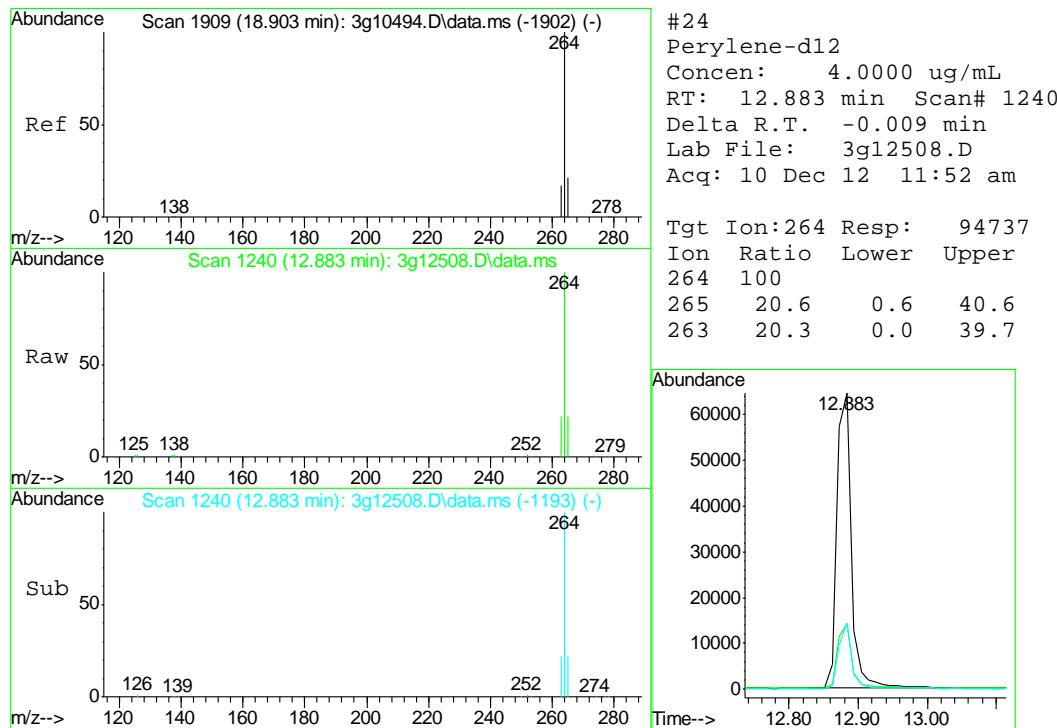
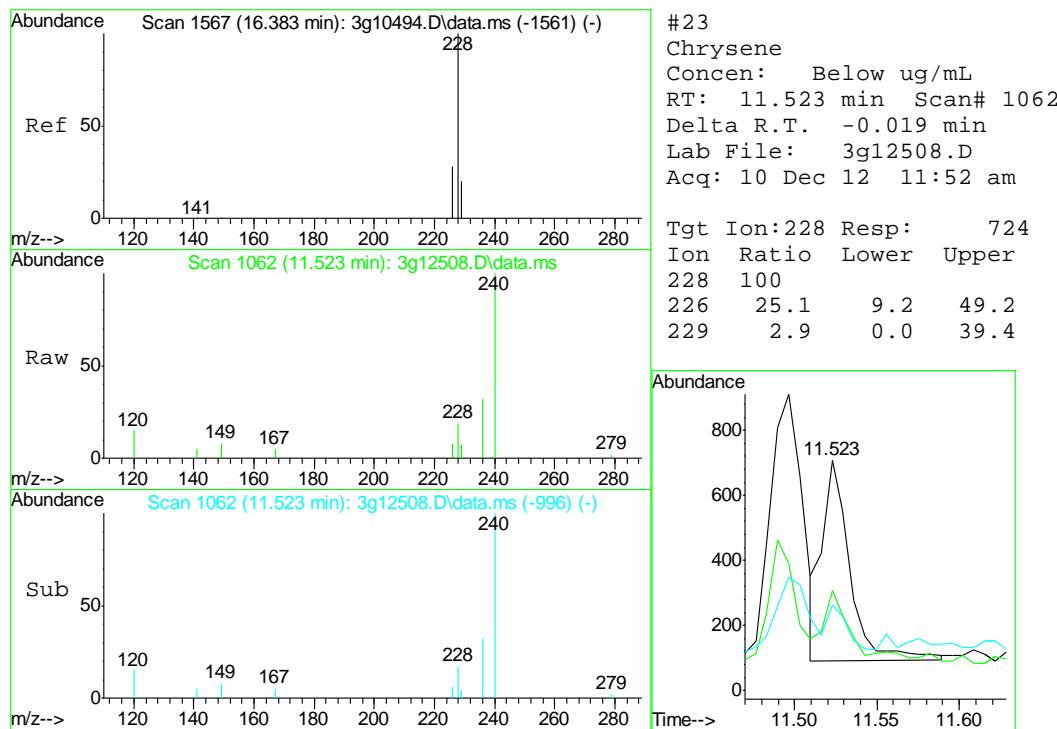


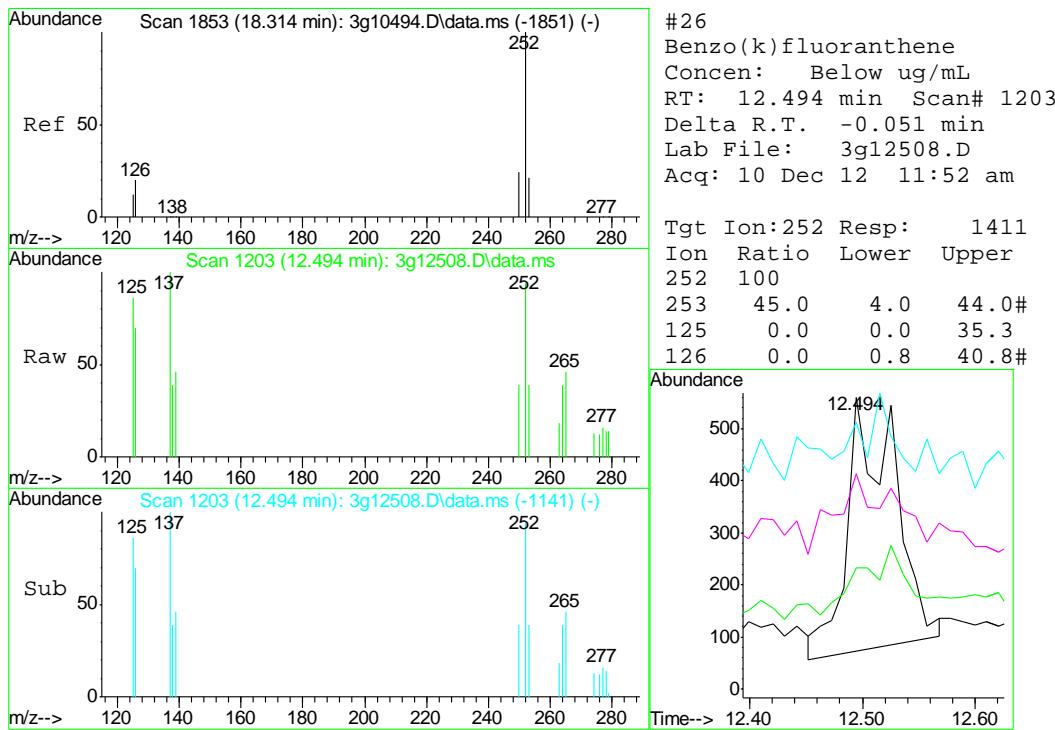
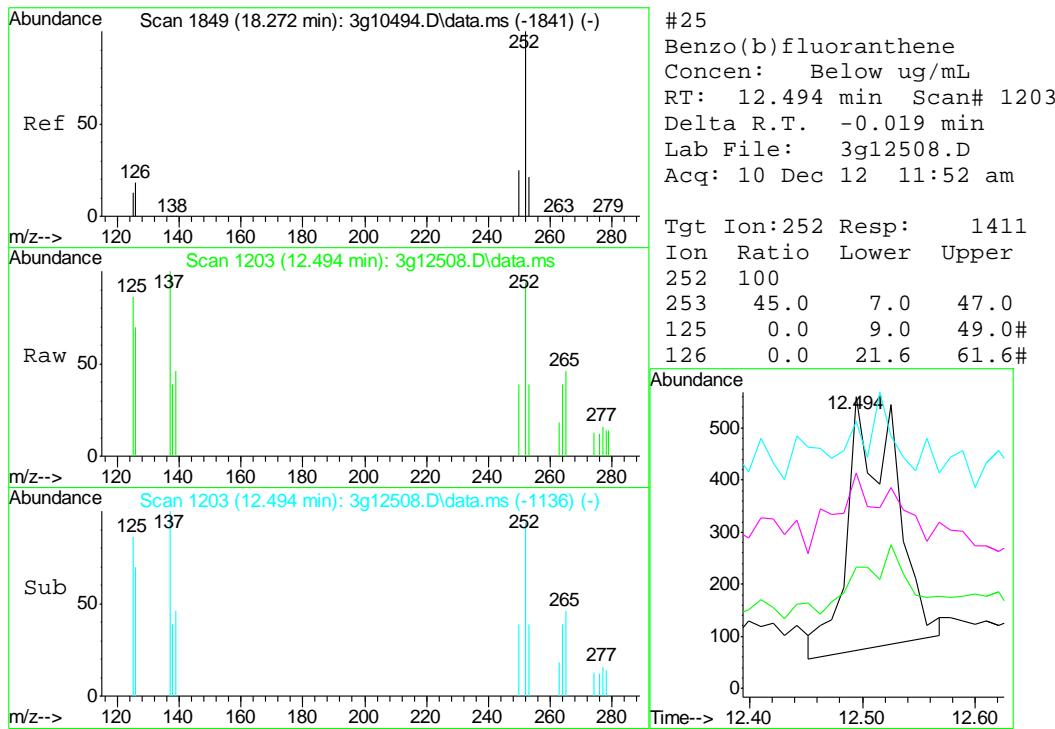


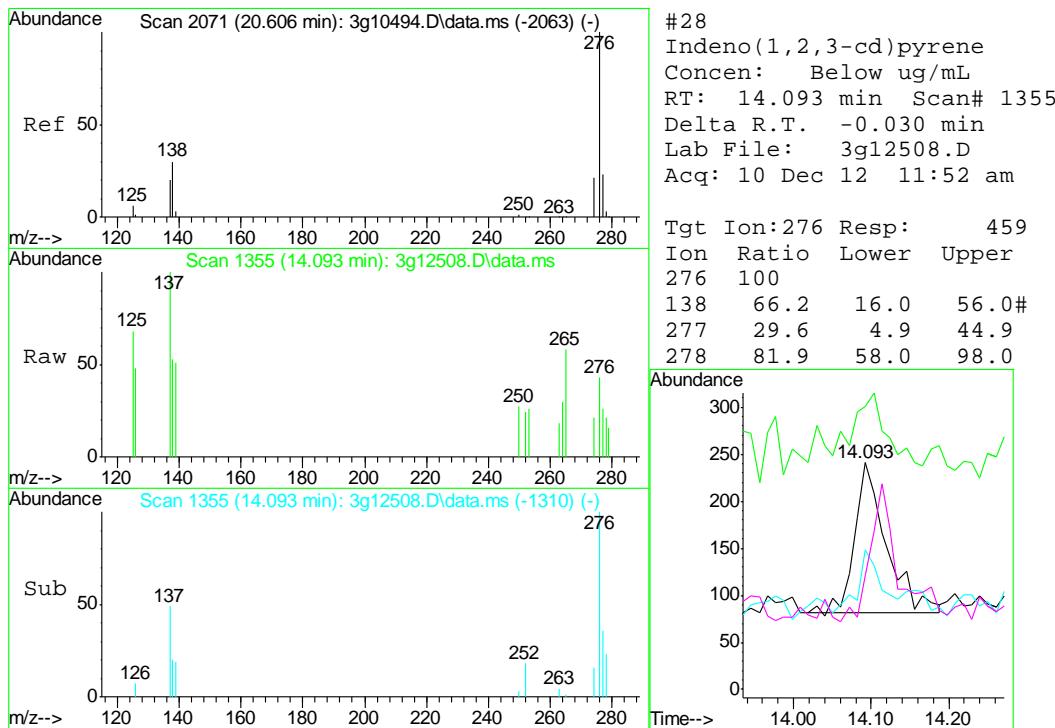
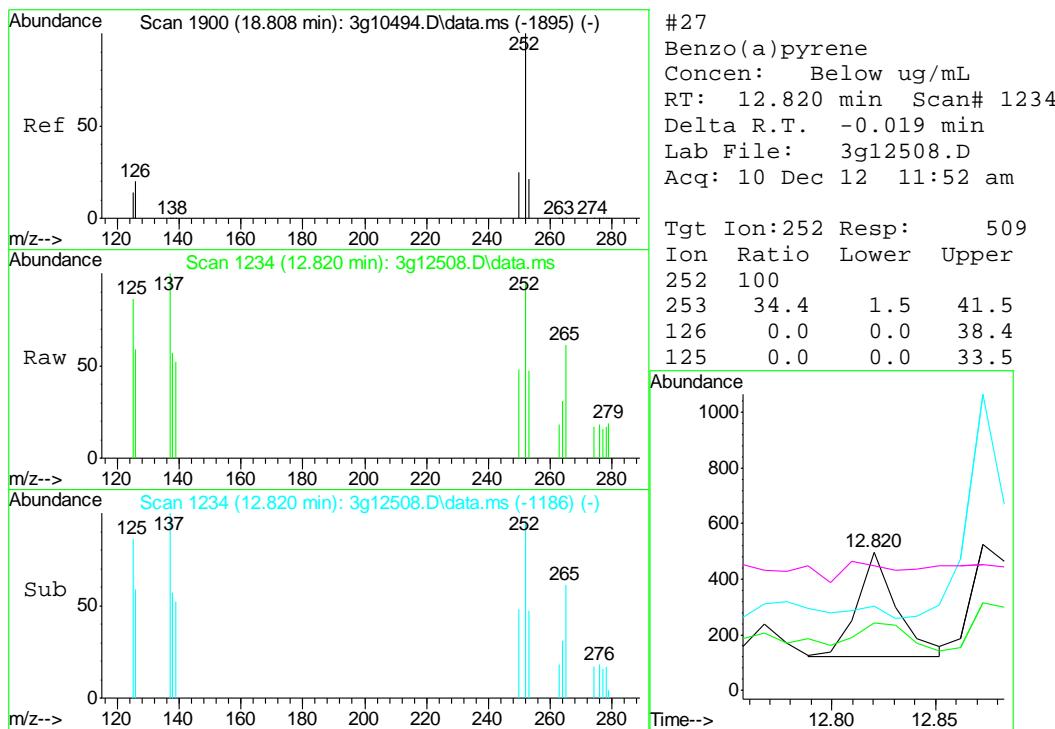


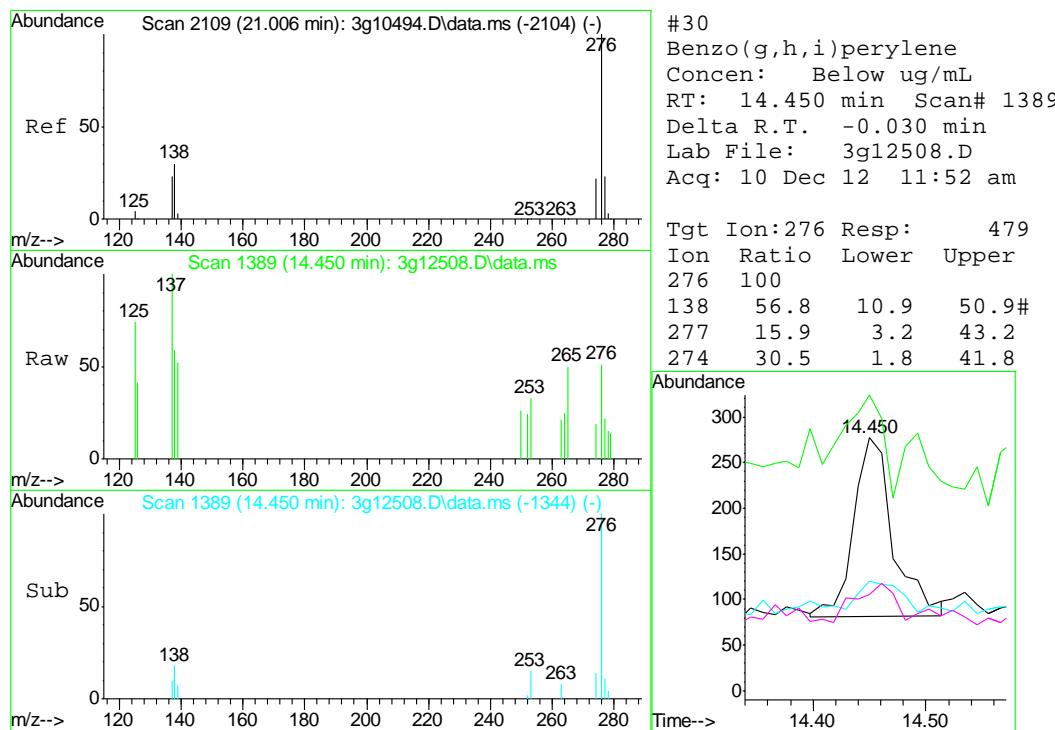
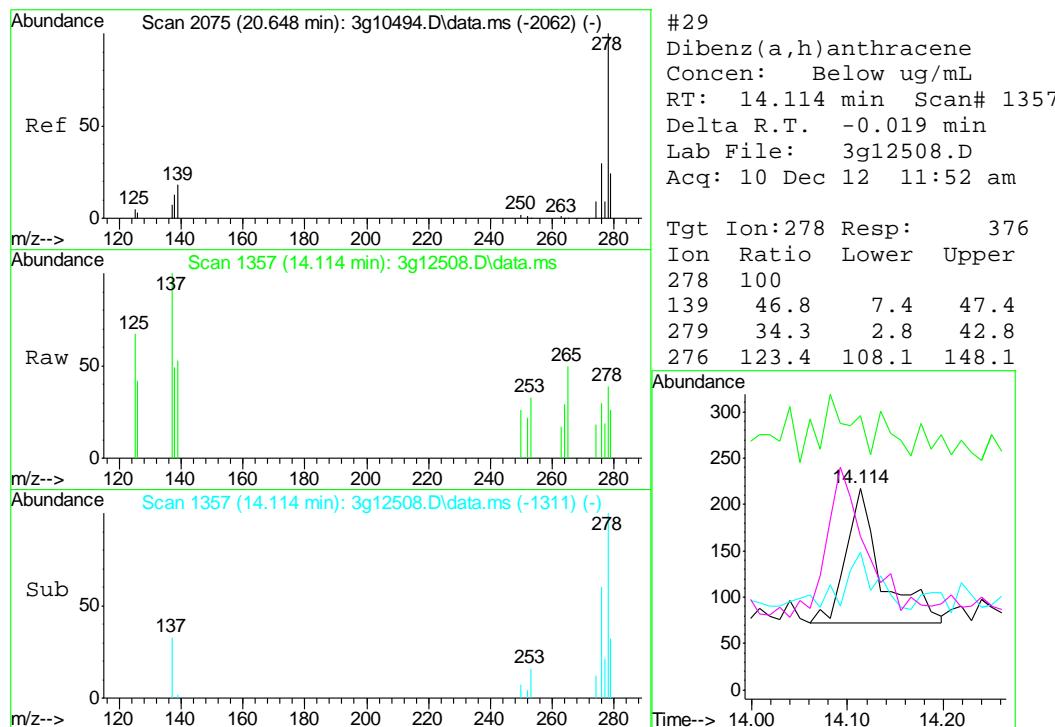














GC Volatiles

QC Data Summaries

Includes the following where applicable:

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

Method Blank Summary

Job Number: D41448
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

D41448-1

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: D41448

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

D41448-1

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

* = Outside of Control Limits.

Matrix Spike/Matrix Spike Duplicate Summary

Page 1 of 1

Job Number: D41448

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

D41448-1

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

Raw Data

Judy Nelson
 12/07/12 12:34

Quantitation Report (QT Reviewed)

Signal #1 : Y:\1\DATA\120612\GB18767.D\FID1A.CH Vial: 8
 Signal #2 : Y:\1\DATA\120612\GB18767.D\FID2B.CH
 Acq On : 6 Dec 2012 3:11 pm Operator: StephK
 Sample : D41448-1, 50X Inst : GC/MS Ins
 Misc : GC3278,GGB1023,,5.015,,100,5,1 Multiplr: 1.00
 IntFile Signal #1: TVH1.E IntFile Signal #2: FB2.E
 Quant Time: Dec 06 17:10:52 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	2611576	83.346 %	m
10) S	1,2,4-Trichlorobenzene (P)	14.37	15514883	95.460 %	

Target Compounds

1) H	TVH-Gasoline	7.23	9017518	0.110 mg/L
4) T	Methyl-t-butyl-ether	0.00	0	N.D. ug/L d
5) T	Benzene	4.14	416527	1.034 ug/L
6) T	Toluene	7.66	1205755	3.043 ug/L
7) T	Ethylbenzene	10.29	166127	0.491 ug/L
8) T	m,p-Xylene	10.47	1268961	3.103 ug/L
9) T	o-Xylene	10.97	228489	0.696 ug/L
11) T	Naphthalene	14.56	3602469	18.258 ug/L

11.11

11

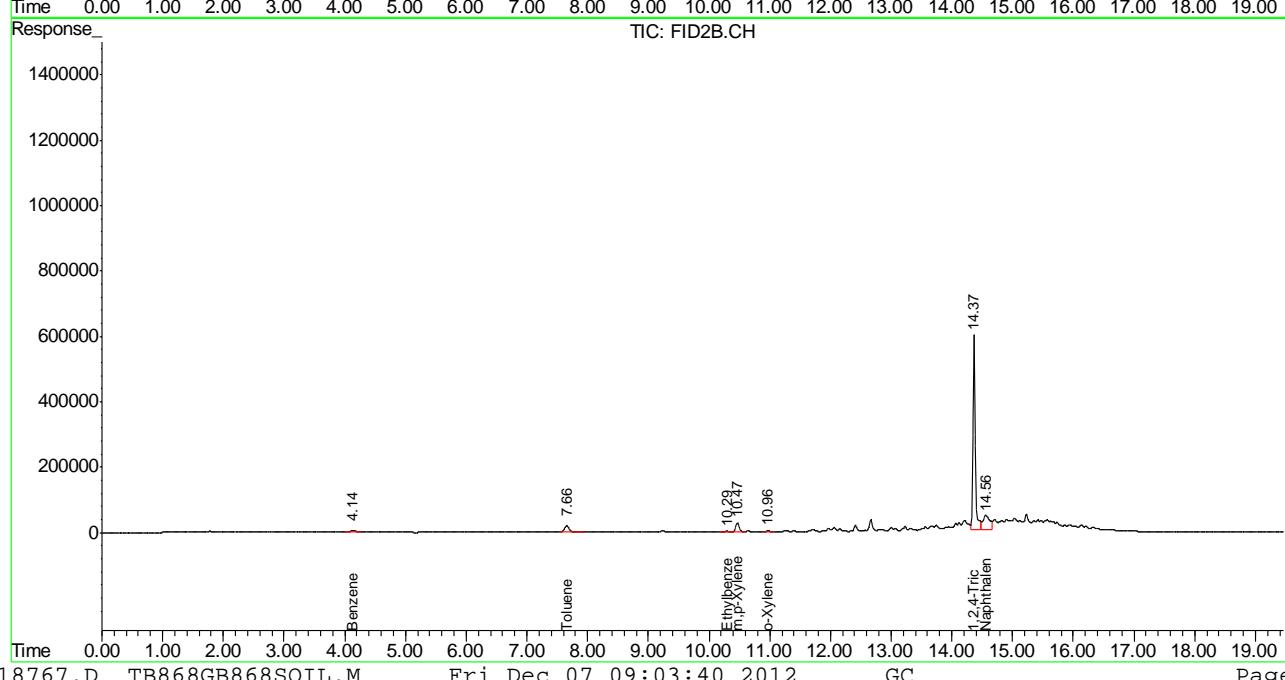
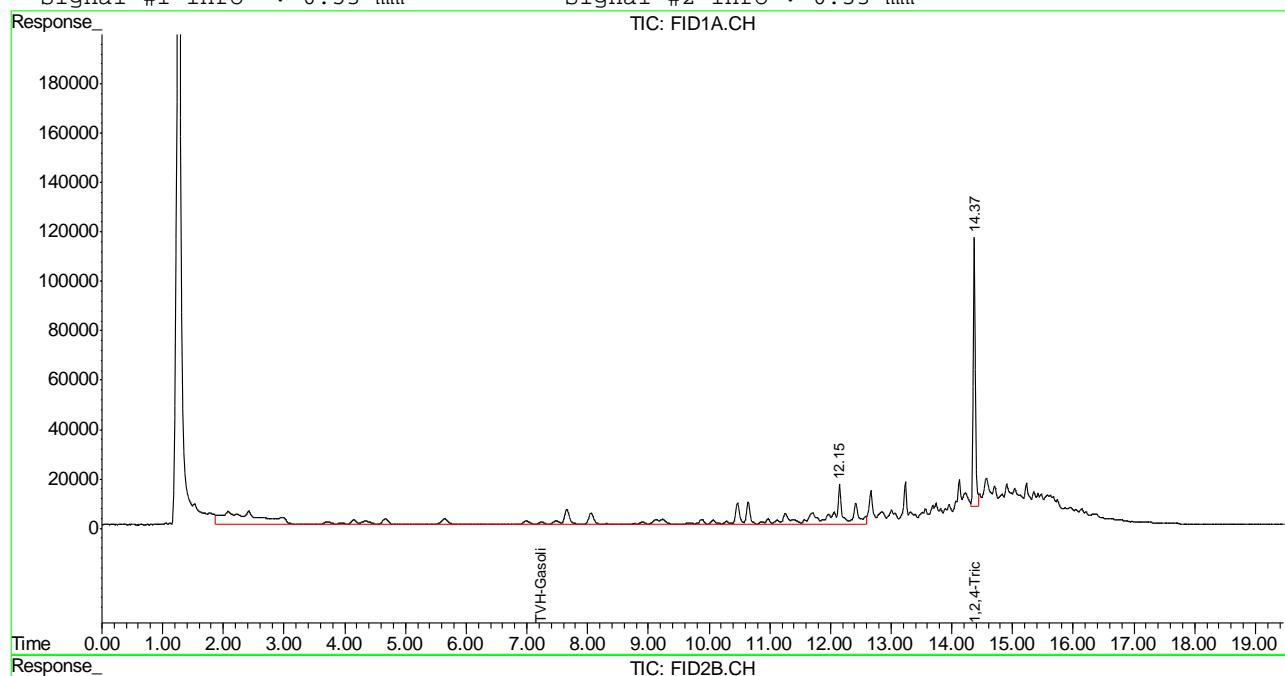
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 GB18767.D TB868GB868SOIL.M Fri Dec 07 09:03:40 2012 GC

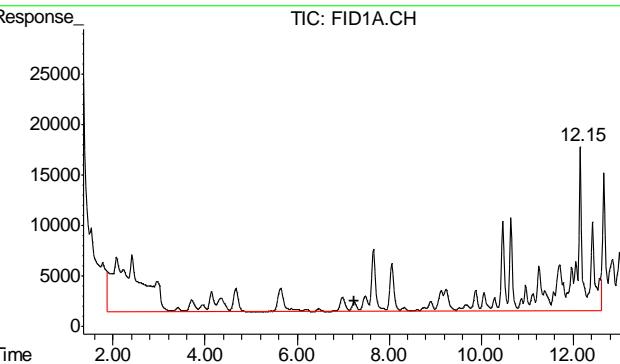
Quantitation Report (QT Reviewed)

Signal #1 : Y:\1\DATA\120612\GB18767.D\FID1A.CH Vial: 8
 Signal #2 : Y:\1\DATA\120612\GB18767.D\FID2B.CH
 Acq On : 6 Dec 2012 3:11 pm Operator: StephK
 Sample : D41448-1, 50X Inst : GC/MS Ins
 Misc : GC3278,GGB1023,5.015,,100,5,1 Multiplr: 1.00
 IntFile Signal #1: TVH1.E IntFile Signal #2: FB2.E
 Quant Time: Dec 6 17:11 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

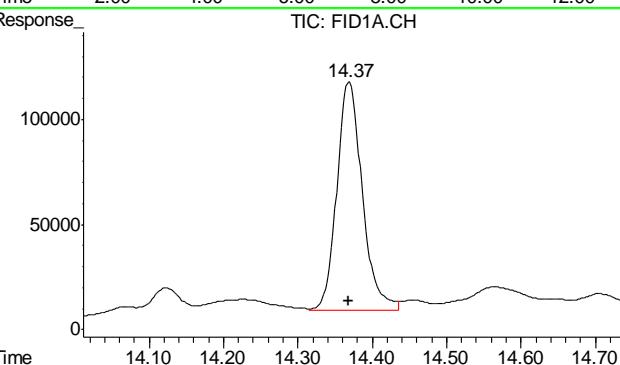
Volume Inj. :
 Signal #1 Phase : DB-624 Signal #2 Phase: DB-624
 Signal #1 Info : 0.53 mm Signal #2 Info : 0.53 mm





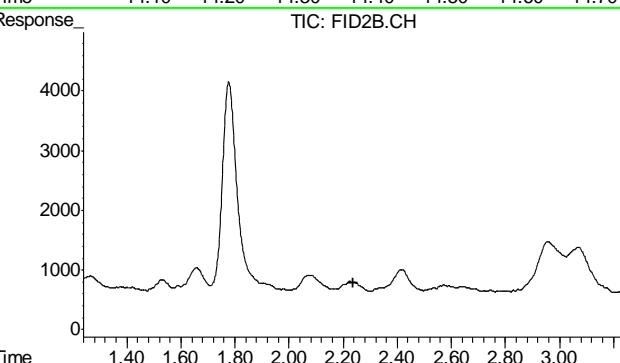
#1 TVH-Gasoline

R.T.: 7.230 min
Delta R.T.: 0.000 min
Response: 9017518
Conc: 0.11 mg/L m



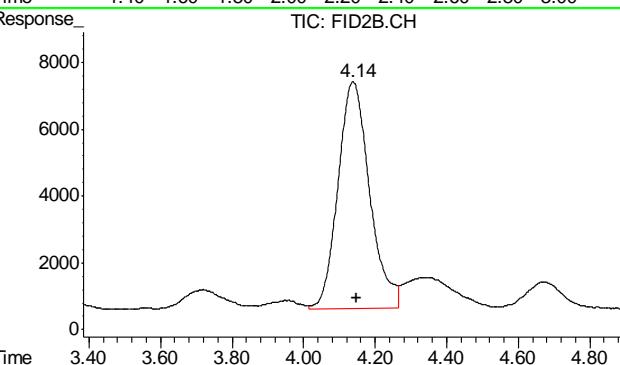
#2 1,2,4-Trichlorobenzene

R.T.: 14.368 min
Delta R.T.: 0.000 min
Response: 2611576
Conc: 83.35 % m



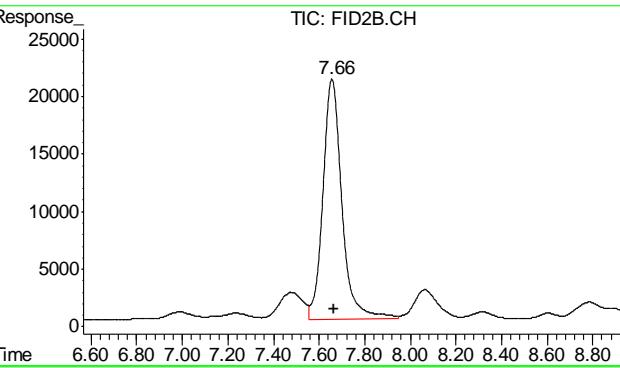
#4 Methyl-t-butyl-ether

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

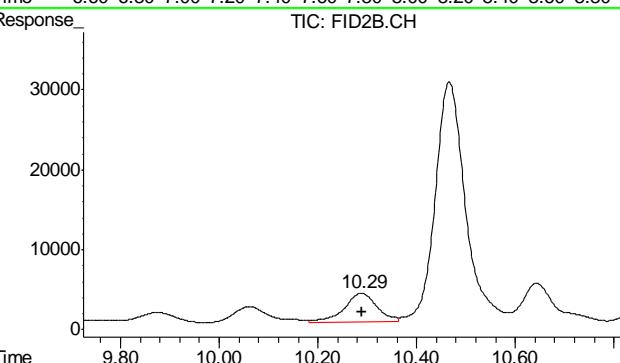


#5 Benzene

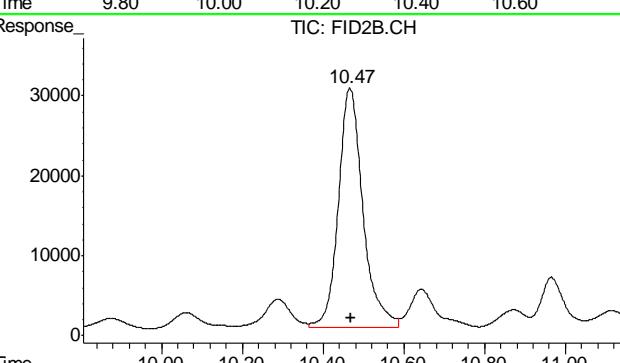
R.T.: 4.138 min
Delta R.T.: -0.009 min
Response: 416527
Conc: 1.03 ug/L



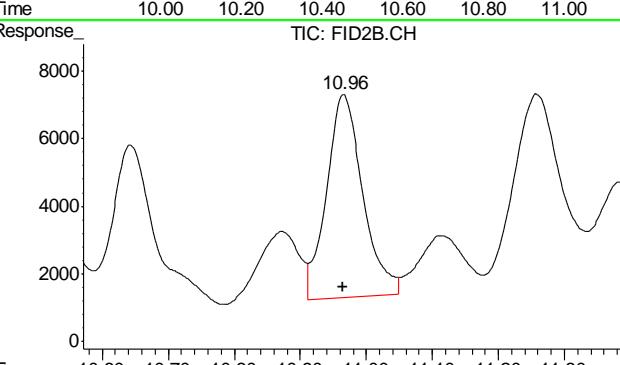
#6 Toluene
R.T.: 7.656 min
Delta R.T.: -0.008 min
Response: 1205755
Conc: 3.04 ug/L



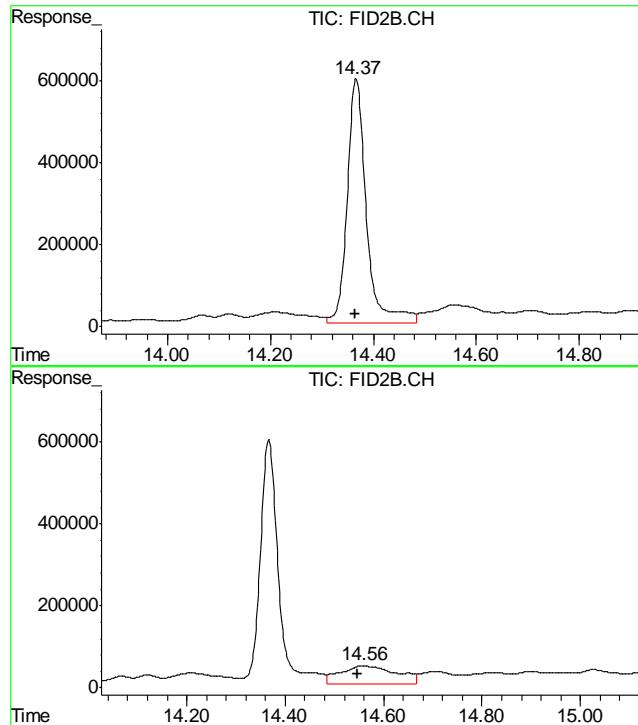
#7 Ethylbenzene
R.T.: 10.289 min
Delta R.T.: 0.000 min
Response: 166127
Conc: 0.49 ug/L



#8 m,p-Xylene
R.T.: 10.466 min
Delta R.T.: -0.003 min
Response: 1268961
Conc: 3.10 ug/L



#9 o-Xylene
R.T.: 10.965 min
Delta R.T.: 0.000 min
Response: 228489
Conc: 0.70 ug/L



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

R.T.: 14.367 min
Delta R.T.: 0.001 min
Response: 15514883
Conc: 95.46 %

#11 Naphthalene

R.T.: 14.558 min
Delta R.T.: 0.010 min
Response: 3602469
Conc: 18.26 ug/L

11.11

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

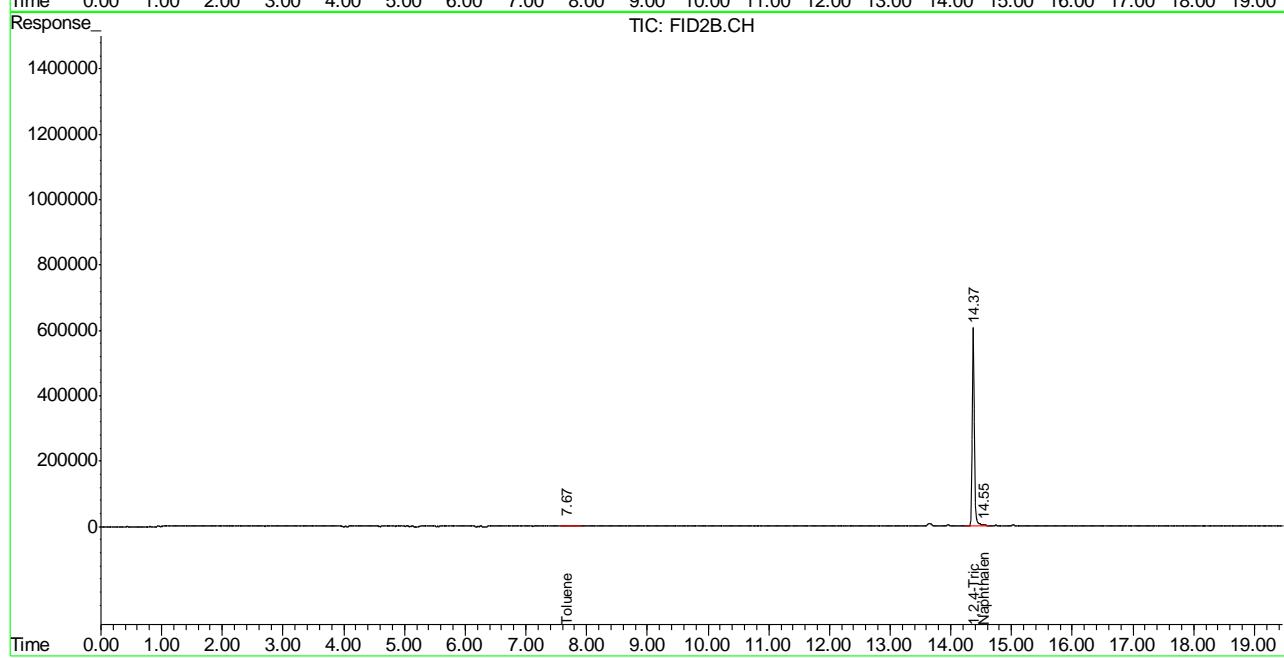
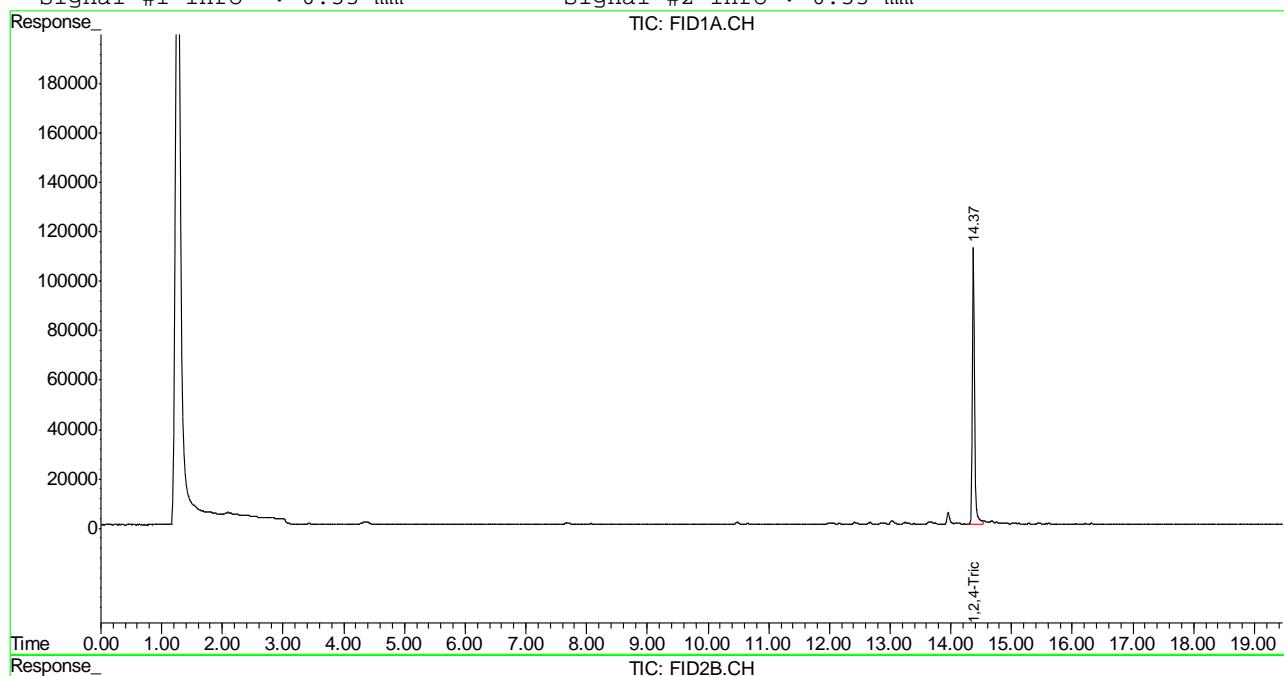
(f)=RT Delta > 1/2 Window (m)=manual int.
 GB18762.D TB868GB868SOIL.M Fri Dec 07 09:03:25 2012 GC

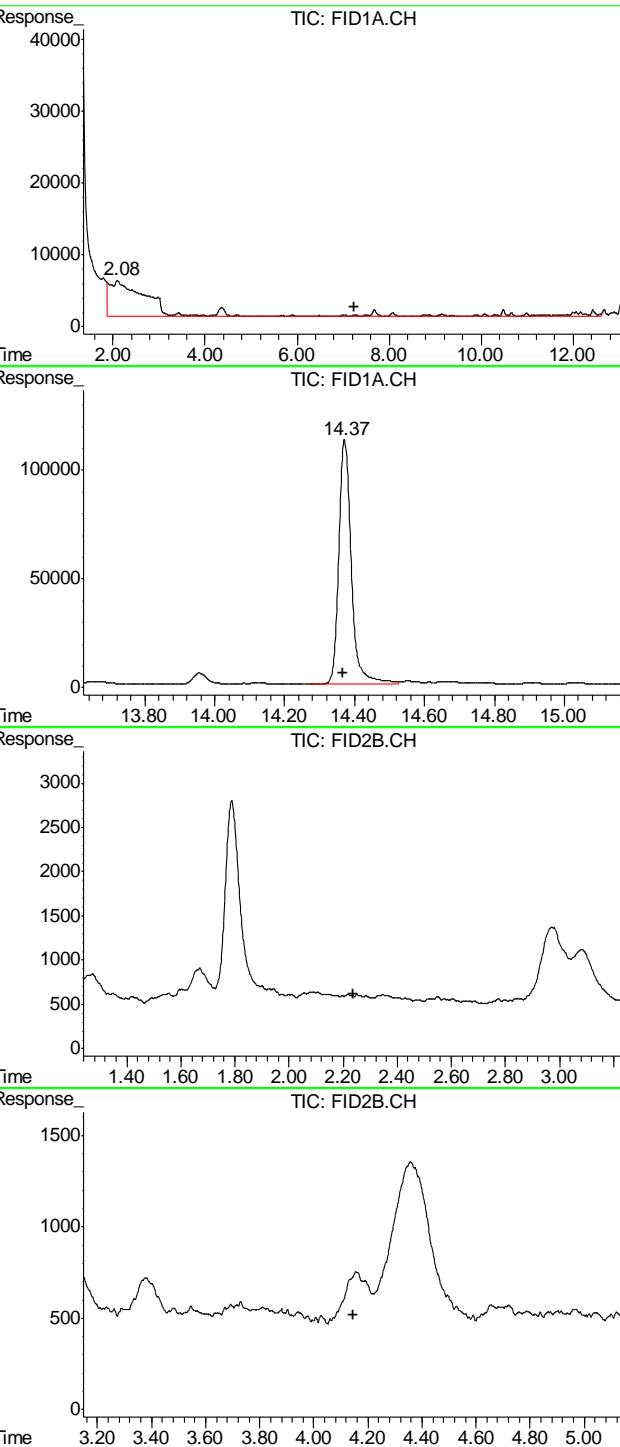
Quantitation Report (QT Reviewed)

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



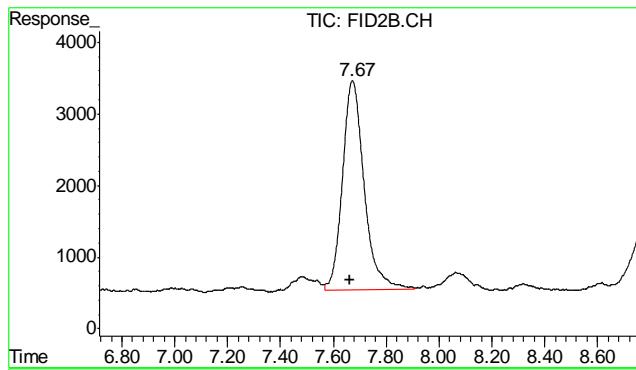


#1 TVH-Gasoline
 R.T.: 7.230 min
 Delta R.T.: 0.000 min
 Response: 3435907
 Conc: N.D.

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

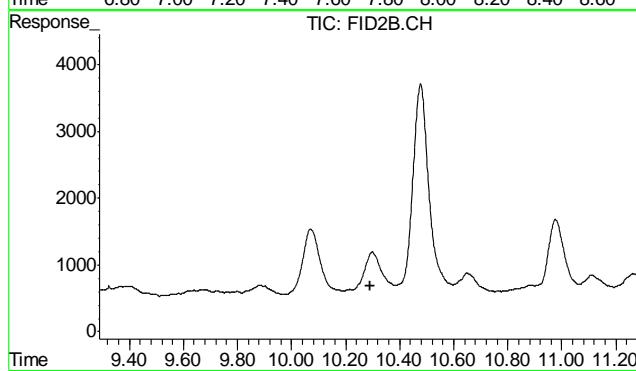
#4 Methyl-t-butyl-ether
 R.T.: 0.000 min
 Exp R.T. : 2.239 min
 Response: 0
 Conc: N.D.

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



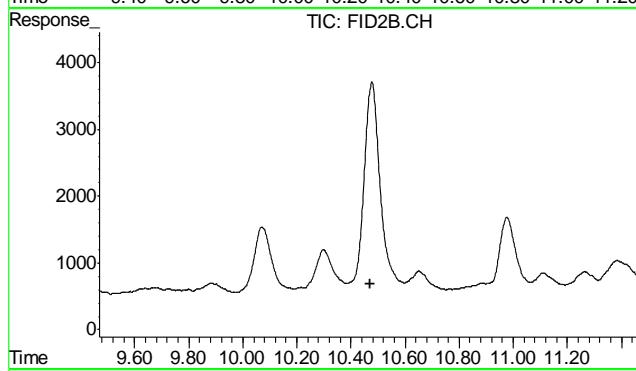
#6 Toluene

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



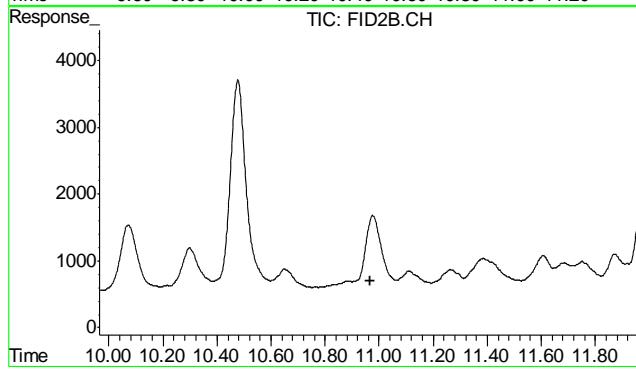
#7 Ethylbenzene

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



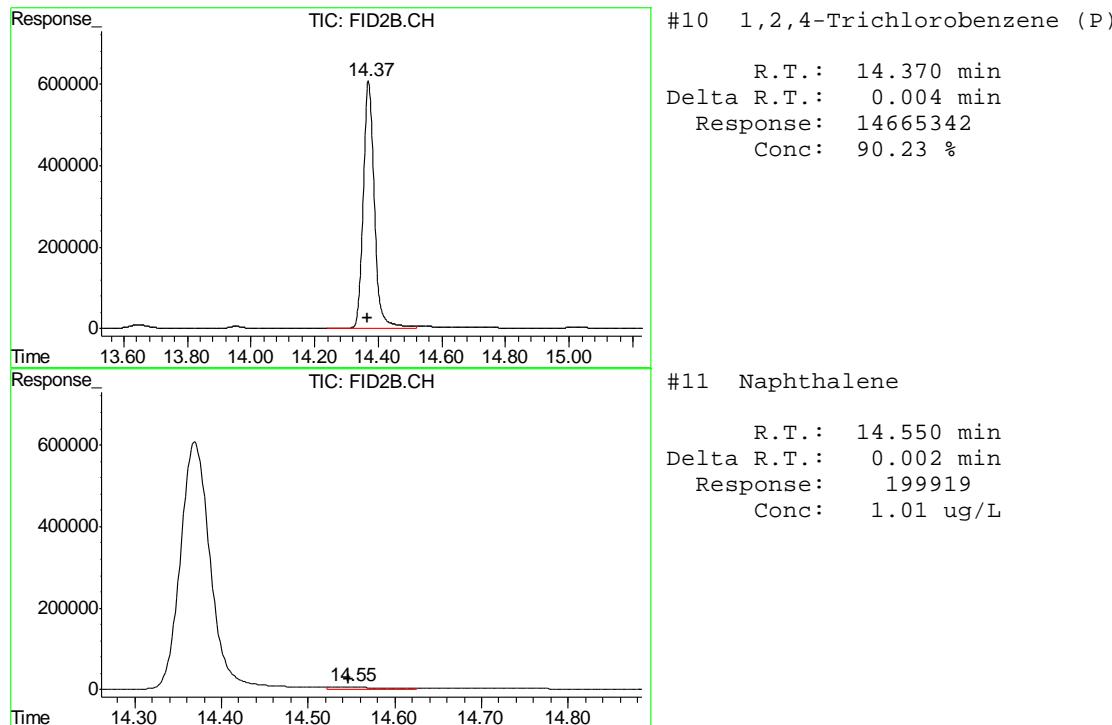
#8 m,p-Xylene

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

QC Data Summaries

Includes the following where applicable:

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

Method Blank Summary

Job Number: D41448
Account: XTOKWR XTO Energy
Project: PCU 296-5A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
OP7053-MB	FH008059.D	1	12/05/12	TR	12/05/12	OP7053	GFH446

The QC reported here applies to the following samples:

Method: SW846-8015B

D41448-1

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: D41448

Account: XTOKWR XTO Energy

Project: PCU 296-5A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
OP7053-BS	FH008060.D	1	12/05/12	TR	12/05/12	OP7053	GFH446

The QC reported here applies to the following samples:

Method: SW846-8015B

D41448-1

CAS No.	Compound	Spike mg/kg	BSP mg/kg	BSP %	Limits
	TPH-DRO (C10-C28)	667	528	79	48-130

CAS No.	Surrogate Recoveries	BSP	Limits
84-15-1	o-Terphenyl	74%	35-130%

* = Outside of Control Limits.

12.2.1
12

Matrix Spike/Matrix Spike Duplicate Summary

Page 1 of 1

Job Number: D41448

Account: XTOKWR XTO Energy

Project: PCU 296-5A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
OP7053-MS	FH008061.D	1	12/05/12	TR	12/05/12	OP7053	GFH446
OP7053-MSD	FH008062.D	1	12/05/12	TR	12/05/12	OP7053	GFH446
D41381-1	FH008063.D	1	12/05/12	TR	12/05/12	OP7053	GFH446

The QC reported here applies to the following samples:

Method: SW846-8015B

D41448-1

CAS No.	Compound	D41381-1		Spike mg/kg	MS mg/kg	MS %	MSD mg/kg	MSD %	RPD	Limits Rec/RPD
		mg/kg	Q							
	TPH-DRO (C10-C28)	18.0		752	534	69	478	61	11	20-168/30
CAS No.	Surrogate Recoveries	MS	MSD	D41381-1		Limits				
84-15-1	o-Terphenyl	56%	54%	68%		35-130%				

* = Outside of Control Limits.

12.3.1
12



GC Semi-volatiles

Raw Data

Manual Integrations
APPROVED
(compounds with "m" flag)

Judy Melson
12/06/12 16:20

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\FH120512\
 Data File : FH008082.D
 Signal(s) : FID1A.ch
 Acq On : 6 Dec 2012 4:13 am
 Operator : TEDR
 Sample : D41448-1
 Misc : OP7053,GFH446,30.07,,,1,1
 ALS Vial : 28 Sample Multiplier: 1

Integration File: autoint1.e
 Quant Time: Dec 06 09:16:55 2012
 Quant Method : C:\msdchem\1\METHODS\YRO-GFH439F.M
 Quant Title : DRO-ORO FRONT
 QLast Update : Fri Nov 30 09:29:08 2012
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. :
 Signal Phase :
 Signal Info :

Compound	R.T.	Response	Conc	Units
<hr/>				
System Monitoring Compounds				
2) s o-Terphenyl	12.737	1557738850	1273.244	ug/mlm
<hr/>				
Target Compounds				
1) H TPH-DRO (C10-C28)	10.422	2900538450	2940.711	ug/ml
<hr/>				

(f)=RT Delta > 1/2 Window

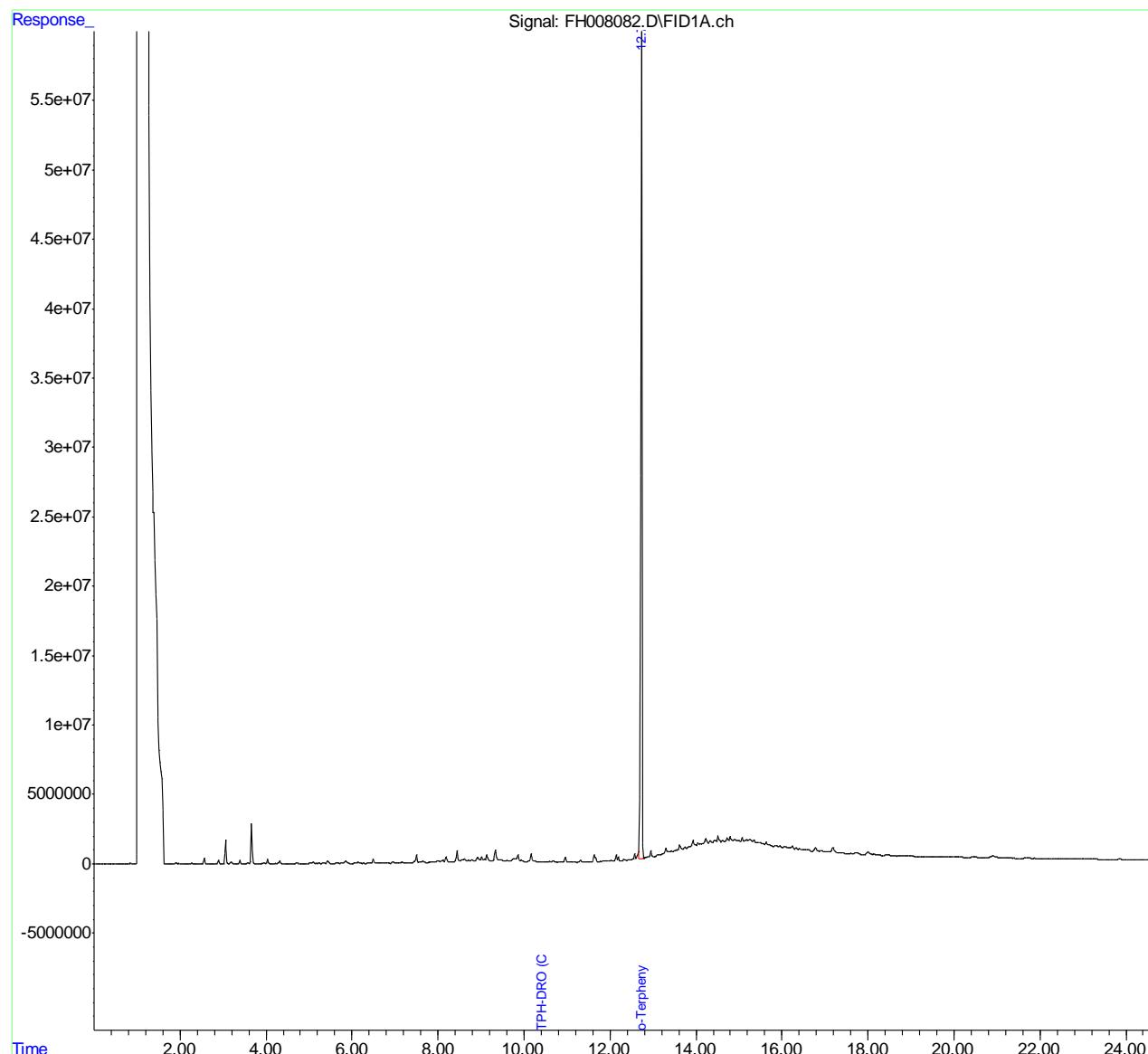
(m)=manual int.

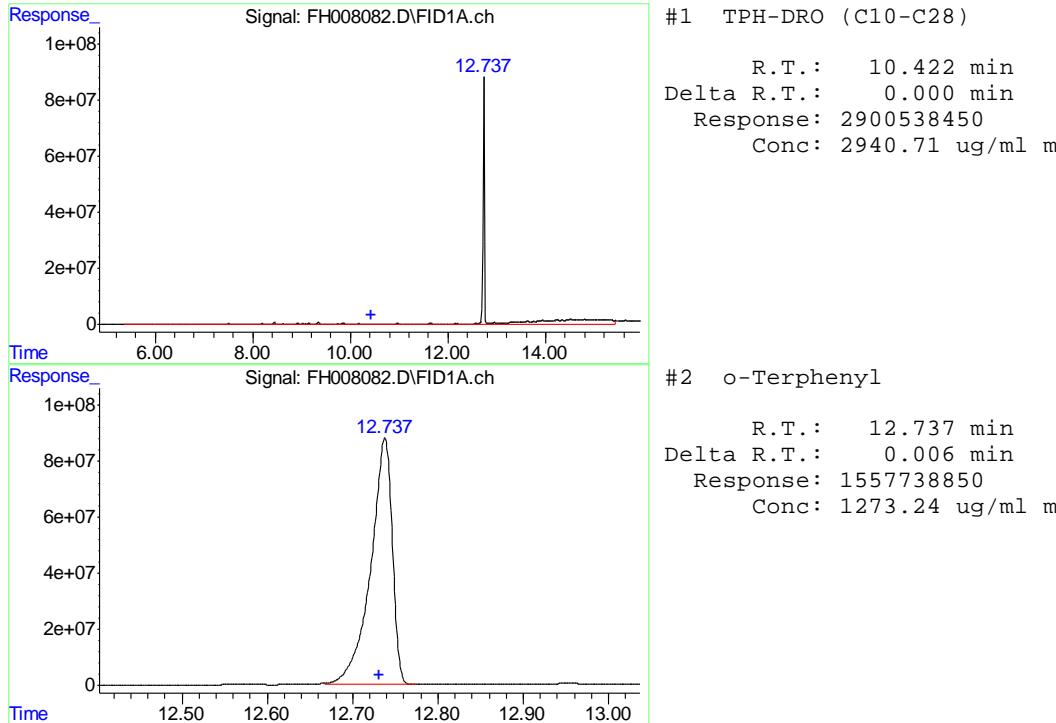
Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\FH120512\
 Data File : FH008082.D
 Signal(s) : FID1A.ch
 Acq On : 6 Dec 2012 4:13 am
 Operator : TEDR
 Sample : D41448-1
 Misc : OP7053,GFH446,30.07,,,1,1
 ALS Vial : 28 Sample Multiplier: 1

Integration File: autoint1.e
 Quant Time: Dec 06 09:16:55 2012
 Quant Method : C:\msdchem\1\METHODS\DRO-GFH439F.M
 Quant Title : DRO-ORO FRONT
 QLast Update : Fri Nov 30 09:29:08 2012
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. :
 Signal Phase :
 Signal Info :





Manual Integrations
APPROVED
(compounds with "m" flag)

Judy Melson
12/06/12 16:19

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\FH120512\
 Data File : FH008059.D
 Signal(s) : FID1A.ch
 Acq On : 5 Dec 2012 2:38 pm
 Operator : TEDR
 Sample : OP7053-MB
 Misc : OP7053,GFH446,30.00,,,1,1
 ALS Vial : 5 Sample Multiplier: 1

Integration File: autoint1.e
 Quant Time: Dec 05 15:04:35 2012
 Quant Method : C:\msdchem\1\METHODS\YRO-GFH439F.M
 Quant Title : DRO-ORO FRONT
 QLast Update : Fri Nov 30 09:29:08 2012
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. :
 Signal Phase :
 Signal Info :

Compound	R.T.	Response	Conc	Units
<hr/>				
System Monitoring Compounds				
2) s o-Terphenyl	12.746	1913998694	1564.439	ug/mlm
<hr/>				
Target Compounds				
1) H TPH-DRO (C10-C28)	10.422	39324729	39.869	ug/ml
<hr/>				

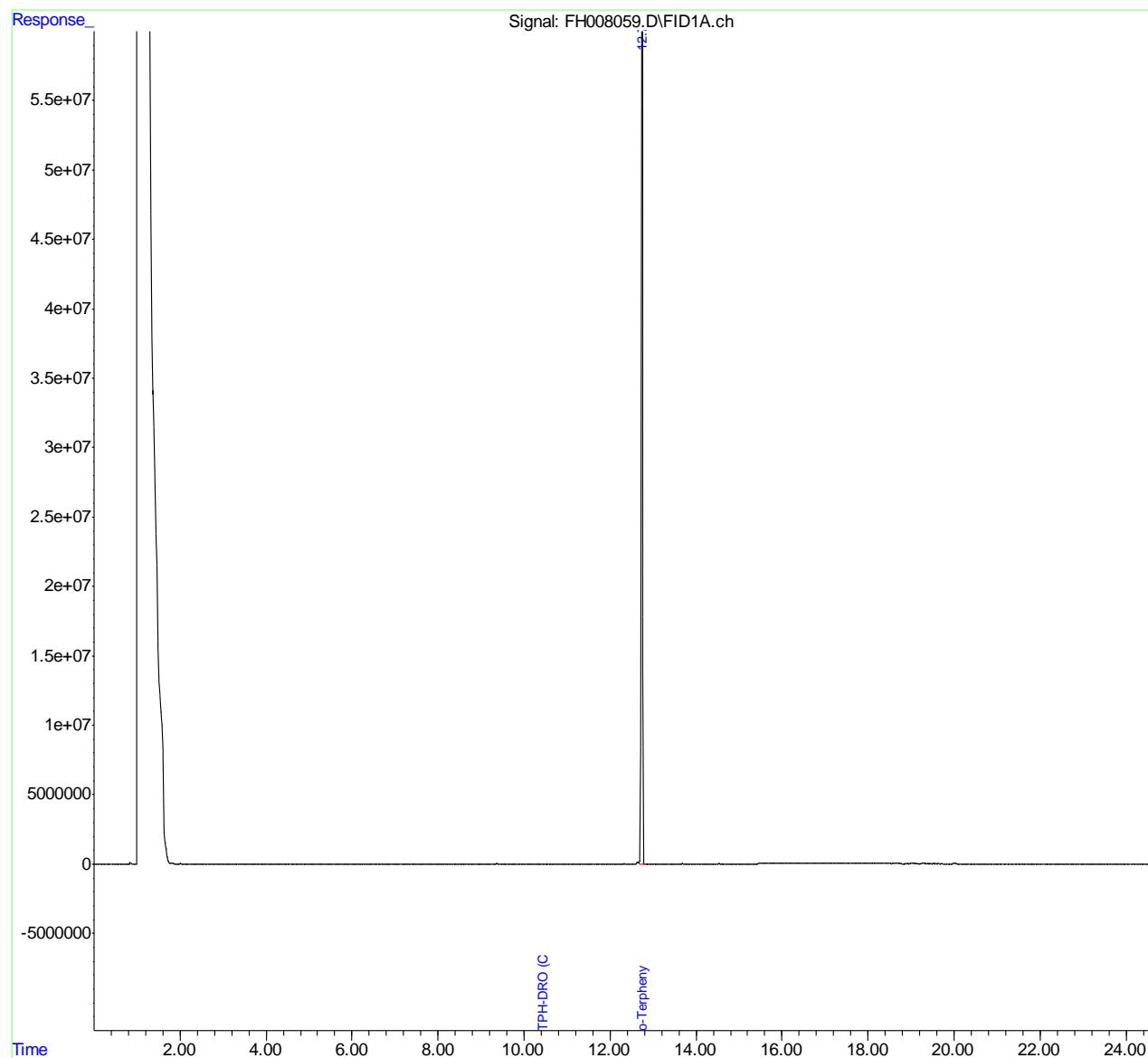
(f)=RT Delta > 1/2 Window (m)=manual int.

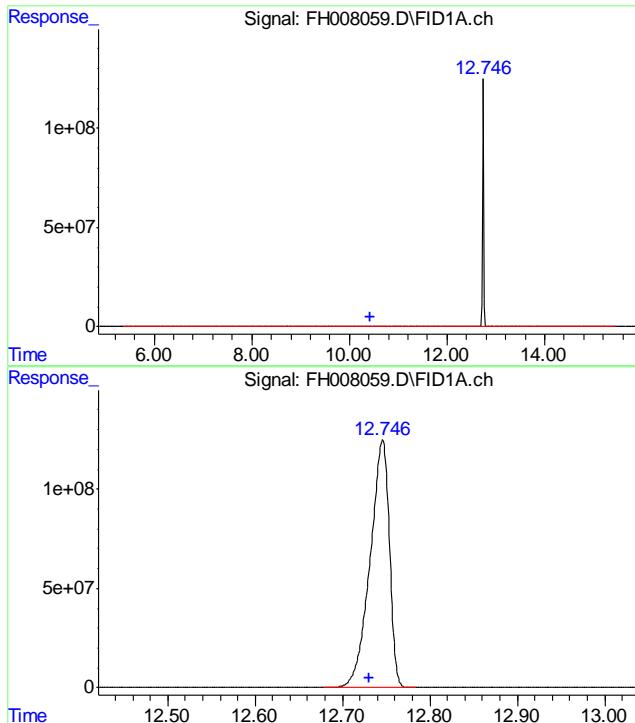
Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\FH120512\
 Data File : FH008059.D
 Signal(s) : FID1A.ch
 Acq On : 5 Dec 2012 2:38 pm
 Operator : TEDR
 Sample : OP7053-MB
 Misc : OP7053,GFH446,30.00,,,1,1
 ALS Vial : 5 Sample Multiplier: 1

Integration File: autoint1.e
 Quant Time: Dec 05 15:04:35 2012
 Quant Method : C:\msdchem\1\METHODS\DRO-GFH439F.M
 Quant Title : DRO-ORO FRONT
 QLast Update : Fri Nov 30 09:29:08 2012
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. :
 Signal Phase :
 Signal Info :





#1 TPH-DRO (C10-C28)

R.T.: 10.422 min
Delta R.T.: 0.000 min
Response: 39324729
Conc: 39.87 ug/ml m

#2 o-Terphenyl

R.T.: 12.746 min
Delta R.T.: 0.015 min
Response: 1913998694
Conc: 1564.44 ug/ml m



Metals Analysis

QC Data Summaries

Includes the following where applicable:

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

BLANK RESULTS SUMMARY
Part 2 - Method Blanks

Login Number: D41448
Account: XTOKRWR - XTO Energy
Project: PCU 296-5A

QC Batch ID: MP9006
Matrix Type: SOLID

Methods: SW846 6010C
Units: mg/kg

Prep Date:

12/05/12

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	0.66	* (a)
Beryllium	1.0	.13	.15		
Boron	5.0	.1	.06		
Cadmium	1.0	.06	.036	0.0	<1.0
Calcium	40	.54	9		
Chromium	1.0	.03	.03	0.080	<1.0
Cobalt	0.50	.04	.07		
Copper	1.0	.12	.15	0.010	<1.0
Iron	7.0	.12	.87		
Lead	5.0	.19	.24	0.070	<5.0
Lithium	0.20	.05	.054		
Magnesium	20	.65	.98		
Manganese	0.50	.12	.022		
Molybdenum	1.0	.21	.08		
Nickel	3.0	.05	.026	0.13	<3.0
Phosphorus	10	1.4	1.9		
Potassium	200	6.1	7		
Selenium	5.0	.48	.36	-0.12	<5.0
Silicon	5.0	.29	.37		
Silver	3.0	.04	.06	-0.070	<3.0
Sodium	40	.59	1.9		
Strontium	5.0	.004	.017		
Thallium	1.0	.29	.53		
Tin	5.0	1.2	2		
Titanium	1.0	.01	.038		
Uranium	5.0	.22	.26		
Vanadium	1.0	.02	.036		
Zinc	3.0	.05	.37	0.21	<3.0

Associated samples MP9006: D41448-1

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

BLANK RESULTS SUMMARY
Part 2 - Method Blanks

Login Number: D41448
Account: XTOKRWR - XTO Energy
Project: PCU 296-5A

QC Batch ID: MP9006
Matrix Type: SOLID

Methods: SW846 6010C
Units: mg/kg

Prep Date:

Metal

(anr) Analyte not requested
(a) All sample results >10x method blank concentration or <RL.

MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D41448
 Account: XTOKRWR - XTO Energy
 Project: PCU 296-5A

QC Batch ID: MP9006
 Matrix Type: SOLID

Methods: SW846 6010C
 Units: mg/kg

Prep Date: 12/05/12

Metal	D41381-1 Original MS	Spikelot ICPALL2	% Rec	QC Limits
Aluminum	anr			
Antimony	anr			
Arsenic	anr			
Barium	228	294	113	58.5N(a) 75-125
Beryllium	anr			
Boron				
Cadmium	0.0	20.9	28.2	73.7N(b) 75-125
Calcium				
Chromium	29.7	46.6	28.2	75.8 75-125
Cobalt	anr			
Copper	15.5	37.5	28.2	80.1 75-125
Iron	anr			
Lead	12.5	53.2	56.4	72.1N(b) 75-125
Lithium				
Magnesium	anr			
Manganese	anr			
Molybdenum	anr			
Nickel	18.9	35.3	28.2	68.8N(b) 75-125
Phosphorus	anr			
Potassium				
Selenium	0.0	45.9	56.4	81.3 75-125
Silicon				
Silver	0.11	9.0	11.3	78.8 75-125
Sodium				
Strontium				
Thallium	anr			
Tin				
Titanium				
Uranium	anr			
Vanadium	anr			
Zinc	41.4	62.9	28.2	76.2 75-125

Associated samples MP9006: D41448-1

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

MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D41448
Account: XTOKRWR - XTO Energy
Project: PCU 296-5A

QC Batch ID: MP9006
Matrix Type: SOLID

Methods: SW846 6010C
Units: mg/kg

Prep Date:

Metal

- (N) Matrix Spike Rec. outside of QC limits
- (anr) Analyte not requested
- (a) Spike recovery indicates possible matrix interference and/or sample nonhomogeneity.
- (b) Spike recovery indicates possible matrix interference.

14.1.2
14

MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D41448
 Account: XTOKRWR - XTO Energy
 Project: PCU 296-5A

QC Batch ID: MP9006
 Matrix Type: SOLID

Methods: SW846 6010C
 Units: mg/kg

Prep Date:

12/05/12

Metal	D41381-1 Original	MSD	Spikelot ICPALL2	% Rec	MSD RPD	QC Limit
Aluminum	anr					
Antimony	anr					
Arsenic	anr					
Barium	228	336	113	95.7	13.3	20
Beryllium	anr					
Boron						
Cadmium	0.0	20.8	28.2	73.3N(a)	0.5	20
Calcium						
Chromium	29.7	48.1	28.2	81.2	3.2	20
Cobalt	anr					
Copper	15.5	38.5	28.2	83.6	2.6	20
Iron	anr					
Lead	12.5	54.2	56.4	73.9N(a)	1.9	20
Lithium						
Magnesium	anr					
Manganese	anr					
Molybdenum	anr					
Nickel	18.9	35.5	28.2	69.5N(a)	0.6	20
Phosphorus	anr					
Potassium						
Selenium	0.0	45.6	56.4	80.8	0.7	20
Silicon						
Silver	0.11	9.0	11.3	78.8	0.0	20
Sodium						
Strontium						
Thallium	anr					
Tin						
Titanium						
Uranium	anr					
Vanadium	anr					
Zinc	41.4	65.2	28.2	84.3	3.6	20

Associated samples MP9006: D41448-1

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

MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D41448
Account: XTOKRWR - XTO Energy
Project: PCU 296-5A

QC Batch ID: MP9006
Matrix Type: SOLID

Methods: SW846 6010C
Units: mg/kg

Prep Date:

Metal

- (N) Matrix Spike Rec. outside of QC limits
(anr) Analyte not requested
(a) Spike recovery indicates possible matrix interference.

SPIKE BLANK AND LAB CONTROL SAMPLE SUMMARY

Login Number: D41448
 Account: XTOKRWR - XTO Energy
 Project: PCU 296-5A

QC Batch ID: MP9006
 Matrix Type: SOLID

Methods: SW846 6010C
 Units: mg/kg

Prep Date: 12/05/12

Metal	BSP Result	Spikelot ICPALL2	% Rec	QC Limits
Aluminum	anr			
Antimony	anr			
Arsenic	anr			
Barium	103	100	103.0	80-120
Beryllium	anr			
Boron				
Cadmium	21.9	25	87.6	80-120
Calcium				
Chromium	23.6	25	94.4	80-120
Cobalt	anr			
Copper	22.8	25	91.2	80-120
Iron	anr			
Lead	45.7	50	91.4	80-120
Lithium				
Magnesium	anr			
Manganese	anr			
Molybdenum	anr			
Nickel	22.8	25	91.2	80-120
Phosphorus	anr			
Potassium				
Selenium	48.3	50	96.6	80-120
Silicon				
Silver	9.6	10	96.0	80-120
Sodium				
Strontium				
Thallium	anr			
Tin				
Titanium				
Uranium	anr			
Vanadium	anr			
Zinc	23.4	25	93.6	80-120

Associated samples MP9006: D41448-1

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

SPIKE BLANK AND LAB CONTROL SAMPLE SUMMARY

Login Number: D41448
Account: XTOKRWR - XTO Energy
Project: PCU 296-5A

QC Batch ID: MP9006
Matrix Type: SOLID

Methods: SW846 6010C
Units: mg/kg

Prep Date:

Metal

(anr) Analyte not requested

SERIAL DILUTION RESULTS SUMMARY

Login Number: D41448
 Account: XTOKRWR - XTO Energy
 Project: PCU 296-5A

QC Batch ID: MP9006
 Matrix Type: SOLID

Methods: SW846 6010C
 Units: ug/l

Prep Date: 12/05/12

Metal	D41381-1 Original	SDL 1:5	%DIF	QC Limits
Aluminum	anr			
Antimony	anr			
Arsenic	anr			
Barium	2020	1950	3.7	0-10
Beryllium	anr			
Boron				
Cadmium	1.00	0.00	100.0(a)	0-10
Calcium				
Chromium	223	267	19.6*(b)	0-10
Cobalt	anr			
Copper	132	142	6.9	0-10
Iron	anr			
Lead	111	122	9.7	0-10
Lithium				
Magnesium	anr			
Manganese	anr			
Molybdenum	anr			
Nickel	168	177	25.4*(b)	0-10
Phosphorus	anr			
Potassium				
Selenium	0.00	0.00	NC	0-10
Silicon				
Silver	4.00	0.00	100.0(a)	0-10
Sodium				
Strontium				
Thallium	anr			
Tin				
Titanium				
Uranium	anr			
Vanadium	anr			
Zinc	367	477	30.0*(b)	0-10

Associated samples MP9006: D41448-1

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

SERIAL DILUTION RESULTS SUMMARY

Login Number: D41448
Account: XTOKRWR - XTO Energy
Project: PCU 296-5A

QC Batch ID: MP9006
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.

BLANK RESULTS SUMMARY
Part 2 - Method Blanks

Login Number: D41448
Account: XTOKRWR - XTO Energy
Project: PCU 296-5A

QC Batch ID: MP9007
Matrix Type: SOLID

Methods: SW846 6020A
Units: mg/kg

Prep Date:

12/05/12

Metal	RL	IDL	MDL	MB raw	final
Aluminum	25	.22	.31		
Antimony	0.20	.0018	.0075		
Arsenic	0.10	.006	.06	-0.0047	<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 MP9007: D41448-1

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

14.2.1
14

MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D41448
 Account: XTOKWR - XTO Energy
 Project: PCU 296-5A

QC Batch ID: MP9007
 Matrix Type: SOLID

Methods: SW846 6020A
 Units: mg/kg

Prep Date:

12/05/12

Metal	D41381-1 Original MS	Spikelot ICPALL2	% Rec	QC Limits
Aluminum				
Antimony				
Arsenic	4.8	51.6	56.4	82.9 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 MP9007: D41448-1

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: D41448
 Account: XTOKRWR - XTO Energy
 Project: PCU 296-5A

QC Batch ID: MP9007
 Matrix Type: SOLID

Methods: SW846 6020A
 Units: mg/kg

Prep Date: 12/05/12

Metal	D41381-1 Original	MSD	Spikelot ICPALL2	% Rec	MSD RPD	QC Limit
Aluminum						
Antimony						
Arsenic	4.8	51.7	56.4	83.1	0.2	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 MP9007: D41448-1

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: D41448
 Account: XTOKRWR - XTO Energy
 Project: PCU 296-5A

QC Batch ID: MP9007
 Matrix Type: SOLID

Methods: SW846 6020A
 Units: mg/kg

Prep Date: 12/05/12

Metal	BSP Result	Spikelot ICPALL2	% Rec	QC Limits
Aluminum				
Antimony				
Arsenic	46.7	50	93.4	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 MP9007: D41448-1

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

14.2.3
14

SERIAL DILUTION RESULTS SUMMARY

Login Number: D41448
 Account: XTOKRWR - XTO Energy
 Project: PCU 296-5A

QC Batch ID: MP9007
 Matrix Type: SOLID

Methods: SW846 6020A
 Units: ug/l

Prep Date: 12/05/12

Metal	D41381-1	Original	SDL	5:25 %DIF	QC Limits
-------	----------	----------	-----	-----------	--------------

Aluminum
 Antimony
 Arsenic 42.2 40.8 3.1 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 MP9007: D41448-1

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

14.2.4
14

BLANK RESULTS SUMMARY
Part 2 - Method Blanks

Login Number: D41448
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: D41448-1

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: D41448
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: D41448-1

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: D41448
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 HGWSR1	Spikelot % Rec	MSD RPD	QC Limit
Mercury	0.26	1.0	0.819	90.4	4.1 20

Associated samples MP9008: D41448-1

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: D41448
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	Limits
Mercury	0.21	0.2	105.0	80-120

Associated samples MP9008: D41448-1

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

14.3.3
14

BLANK RESULTS SUMMARY
Part 2 - Method Blanks

Login Number: D41448
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: D41448-1A

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

BLANK RESULTS SUMMARY
Part 2 - Method Blanks

Login Number: D41448
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: D41448
 Account: XTOKWR - 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: D41448-1A

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

MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D41448
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: D41448
 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: D41448-1A

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

MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D41448
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: D41448
 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: D41448-1A

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

SPIKE BLANK AND LAB CONTROL SAMPLE SUMMARY

Login Number: D41448
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

14.4.3
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SERIAL DILUTION RESULTS SUMMARY

Login Number: D41448
 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: D41448-1A

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

SERIAL DILUTION RESULTS SUMMARY

Login Number: D41448
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

QC Data Summaries

Includes the following where applicable:

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

METHOD BLANK AND SPIKE RESULTS SUMMARY
GENERAL CHEMISTRY

Login Number: D41448
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	GN17962			su	8.00	7.99	99.9	99.3-100.7%

Associated Samples:

Batch GP8845: D41448-1

Batch GP8856: D41448-1

Batch GN17962: D41448-1

(*) Outside of QC limits

DUPLICATE RESULTS SUMMARY
GENERAL CHEMISTRY

Login Number: D41448
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: D41448-1
(*) Outside of QC limits

MATRIX SPIKE RESULTS SUMMARY
GENERAL CHEMISTRY

Login Number: D41448
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: D41448-1

(*) Outside of QC limits

(N) Matrix Spike Rec. outside of QC limits

MATRIX SPIKE DUPLICATE RESULTS SUMMARY
GENERAL CHEMISTRY

Login Number: D41448
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: D41448-1

(*) Outside of QC limits

(N) Matrix Spike Rec. outside of QC limits