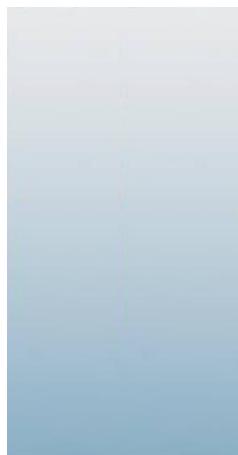




11/27/12



Technical Report for

XTO Energy

PCU 296-5A

1210-04

Accutest Job Number: D41014

Sampling Date: 11/14/12

Report to:

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

Total number of pages in report: 149



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, appearing to read "H. Madadian".

Brad Madadian
Laboratory Director

Client Service contact: Renea Jackson 303-425-6021

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

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

Table of Contents

-1-

Section 1: Sample Summary	4	
Section 2: Case Narrative/Conformance Summary	5	
Section 3: Summary of Hits	8	
Section 4: Sample Results	9	
4.1: D41014-1: CUT 2 SUBLINER (COMP)	10	
4.2: D41014-1A: CUT 2 SUBLINER (COMP)	16	
Section 5: Misc. Forms	18	
5.1: Chain of Custody	19	
Section 6: GC/MS Volatiles - QC Data Summaries	21	
6.1: Method Blank Summary	22	
6.2: Blank Spike Summary	23	
6.3: Matrix Spike/Matrix Spike Duplicate Summary	24	
Section 7: GC/MS Volatiles - Raw Data	25	
7.1: Samples	26	
7.2: Method Blanks	40	
Section 8: GC/MS Semi-volatiles - QC Data Summaries	50	
8.1: Method Blank Summary	51	
8.2: Blank Spike Summary	52	
8.3: Matrix Spike/Matrix Spike Duplicate Summary	53	
Section 9: GC/MS Semi-volatiles - Raw Data	54	
9.1: Samples	55	
9.2: Method Blanks	72	
Section 10: GC Volatiles - QC Data Summaries	89	
10.1: Method Blank Summary	90	
10.2: Blank Spike Summary	91	
10.3: Matrix Spike/Matrix Spike Duplicate Summary	92	
Section 11: GC Volatiles - Raw Data	93	
11.1: Samples	94	
11.2: Method Blanks	99	
Section 12: GC Semi-volatiles - QC Data Summaries	104	
12.1: Method Blank Summary	105	
12.2: Blank Spike Summary	106	
12.3: Matrix Spike/Matrix Spike Duplicate Summary	107	
Section 13: GC Semi-volatiles - Raw Data	108	
13.1: Samples	109	
13.2: Method Blanks	112	
Section 14: Metals Analysis - QC Data Summaries	115	
14.1: Prep QC MP8913: Ba,Cd,Cr,Cu,Pb,Ni,Se,Ag,Zn	116	
14.2: Prep QC MP8914: As	126	
14.3: Prep QC MP8915: Ca,Mg,Na,Sodium Adsorption Ratio	131	
14.4: Prep QC MP8936: Hg	141	
Section 15: General Chemistry - QC Data Summaries	145	

Table of Contents

-2-

15.1: Method Blank and Spike Results Summary	146
15.2: Duplicate Results Summary	147
15.3: Matrix Spike Results Summary	148
15.4: Matrix Spike Duplicate Results Summary	149

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



Sample Summary

XTO Energy

Job No: D41014

PCU 296-5A

Project No: 1210-04

Sample Number	Collected Date	Time By	Received	Matrix Code	Type	Client Sample ID
D41014-1	11/14/12	13:45 DS	11/16/12	SO	Soil	CUT 2 SUBLINER (COMP)
D41014-1A	11/14/12	13:45 DS	11/16/12	SO	Soil	CUT 2 SUBLINER (COMP)

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



CASE NARRATIVE / CONFORMANCE SUMMARY

Client: XTO Energy

Job No D41014

Site: PCU 296-5A

Report Date 11/27/2012 2:15:27 PM

On 11/16/2012, 1 sample(s), 0 Trip Blank(s), and 0 Field Blank(s) were received at Accutest Mountain States (AMS) at a temperature of 2.7 °C. The samples were intact and properly preserved, unless noted below. An AMS Job Number of D41014 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: V5V1506
------------------	--------------------------

- All samples were analyzed within the recommended method holding time.
- All method blanks for this batch meet method specific criteria.
- Sample(s) D41036-1MS, D41036-1MSD were used as the QC samples indicated.

Extractables by GCMS By Method SW846 8270C BY SIM

Matrix SO	Batch ID: OP6988
------------------	-------------------------

- All samples were extracted within the recommended method holding time.
- All samples were analyzed within the recommended method holding time.
- Sample(s) D41014-1MS, D41014-1MSD 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: GGB1010
------------------	--------------------------

- All samples were analyzed within the recommended method holding time.
- All method blanks for this batch meet method specific criteria.
- Sample(s) D41015-1MS, D41015-1MSD were used as the QC samples indicated.

Extractables by GC By Method SW846-8015B

Matrix SO	Batch ID: OP6979
------------------	-------------------------

- All samples were extracted 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) D41044-9MS, D41044-9MSD were used as the QC samples indicated.

Metals By Method SW846 6010C

Matrix AQ

Batch ID: MP8915

- 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) D41042-1AMS, D41042-1AMSD, D41042-1ASDL were used as the QC samples for the metals analysis.
- The serial dilution RPD(s) for Magnesium, Sodium are outside control limits for sample MP8915-SD1. Probable cause due to sample homogeneity.
- MP8915-SD1 for Magnesium: Serial dilution indicates possible matrix interference.
- MP8915-SD1 for Sodium: Serial dilution indicates possible matrix interference.

Matrix SO

Batch ID: MP8913

- 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) D41013-1MS, D41013-1MSD, D41013-1SDL were used as the QC samples for the metals analysis.
- The matrix spike (MS) recovery(s) of Nickel, Silver are outside control limits. Spike recovery indicates possible matrix interference.
- The matrix spike duplicate (MSD) recovery(s) of Chromium, Nickel, Silver are outside control limits. Probable cause due to matrix interference.
- The matrix spike (MS) recovery(s) of Barium are outside control limits. Spike amount low relative to the sample amount. Refer to lab control or spike blank for recovery information.
- The serial dilution RPD(s) for Selenium, Silver, Barium, Chromium, Nickel, Zinc are outside control limits for sample MP8913-SD1. Percent difference acceptable due to low initial sample concentration (< 50 times IDL).
- MP8913-SD1 for Nickel: Serial dilution indicates possible matrix interference.
- MP8913-SD1 for Chromium: Serial dilution indicates possible matrix interference.
- MP8913-SD1 for Barium: Serial dilution indicates possible matrix interference.
- MP8913-SD1 for Zinc: Serial dilution indicates possible matrix interference.

Metals By Method SW846 6020A

Matrix SO

Batch ID: MP8914

- 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) D41013-1MS, D41013-1MSD, D41013-1SDL were used as the QC samples for the metals analysis.

Metals By Method SW846 7471B

Matrix SO

Batch ID: MP8936

- 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) D40988-1MS, D40988-1MSD were used as the QC samples for the metals analysis.

Wet Chemistry By Method ASTM D1498-76M

Matrix SO

Batch ID: GN17722

- Sample(s) D41014-1DUP were used as the QC samples for the Redox Potential Vs H₂ analysis.

Wet Chemistry By Method SM19 2540B M

Matrix SO

Batch ID: GN17724

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

Wet Chemistry By Method SW846 3060/7196A M

Matrix SO

Batch ID: R15202

- The data for SW846 3060/7196A M meets quality control requirements.
- D41014-1 for Chromium, Trivalent: Calculated as: (Chromium) - (Chromium, Hexavalent)

Wet Chemistry By Method SW846 3060A/7196A

Matrix SO

Batch ID: GP8709

- 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) D40985-1DUP, D40985-1MS, D40985-1MSD were used as the QC samples for the Chromium, Hexavalent analysis.

Wet Chemistry By Method SW846 9045D

Matrix SO

Batch ID: GN17719

- The following samples were run outside of holding time for method SW846 9045D: D41014-1

Wet Chemistry By Method USDA HANDBOOK 60

Matrix SO

Batch ID: MP8915

- D41014-1A 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: D41014
Account: XTO Energy
Project: PCU 296-5A
Collected: 11/14/12

3

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

D41014-1 CUT 2 SUBLINER (COMP)

Benzene	0.0444 J	0.065	0.032	mg/kg	SW846 8260B
Toluene	0.158	0.13	0.065	mg/kg	SW846 8260B
Ethylbenzene	0.0330 J	0.13	0.025	mg/kg	SW846 8260B
Xylene (total)	0.161 J	0.26	0.13	mg/kg	SW846 8260B
Naphthalene	0.0348	0.013	0.012	mg/kg	SW846 8270C BY SIM
TPH-DRO (C10-C28)	24.7	15	10	mg/kg	SW846-8015B
Arsenic	13.0	0.11		mg/kg	SW846 6020A
Barium	1710	1.1		mg/kg	SW846 6010C
Chromium	65.4	1.1		mg/kg	SW846 6010C
Copper	8.1	1.1		mg/kg	SW846 6010C
Lead	9.1	5.7		mg/kg	SW846 6010C
Nickel	20.5	3.4		mg/kg	SW846 6010C
Zinc	38.2	3.4		mg/kg	SW846 6010C
Specific Conductivity	353	1.0		umhos/cm	SM2510B-1997 MOD
Chromium, Trivalent ^a	65.4	2.1		mg/kg	SW846 3060/7196A M
Redox Potential Vs H2	94.2			mv	ASTM D1498-76M
pH	9.90			su	SW846 9045D

D41014-1A CUT 2 SUBLINER (COMP)

Calcium	14.0	2.0	mg/l	SW846 6010C
Magnesium	4.53	1.0	mg/l	SW846 6010C
Sodium	61.8	2.0	mg/l	SW846 6010C
Sodium Adsorption Ratio ^b	3.67		ratio	USDA HANDBOOK 60

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

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



4

Sample Results

Report of Analysis

Accutest Laboratories

Report of Analysis

Page 1 of 1

Client Sample ID: CUT 2 SUBLINER (COMP)**Lab Sample ID:** D41014-1**Matrix:** SO - Soil**Method:** SW846 8260B**Project:** PCU 296-5A**Date Sampled:** 11/14/12**Date Received:** 11/16/12**Percent Solids:** 86.4

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	5V24680.D	1	11/19/12	BD	n/a	n/a	V5V1506
Run #2							

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

Purgeable Aromatics

CAS No.	Compound	Result	RL	MDL	Units	Q
71-43-2	Benzene	0.0444	0.065	0.032	mg/kg	J
108-88-3	Toluene	0.158	0.13	0.065	mg/kg	
100-41-4	Ethylbenzene	0.0330	0.13	0.025	mg/kg	J
1330-20-7	Xylene (total)	0.161	0.26	0.13	mg/kg	J

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
2037-26-5	Toluene-D8	101%		64-130%
460-00-4	4-Bromofluorobenzene	100%		62-131%
17060-07-0	1,2-Dichloroethane-D4	96%		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:	CUT 2 SUBLINER (COMP)	Date Sampled:	11/14/12
Lab Sample ID:	D41014-1	Date Received:	11/16/12
Matrix:	SO - Soil	Percent Solids:	86.4
Method:	SW846 8270C BY SIM	SW846 3546	
Project:	PCU 296-5A		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	3G12231.D	1	11/26/12	SM	11/20/12	OP6988	E3G577
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	0.0096	0.0050	mg/kg	
120-12-7	Anthracene	ND	0.0096	0.0050	mg/kg	
56-55-3	Benzo(a)anthracene	ND	0.0096	0.0050	mg/kg	
205-99-2	Benzo(b)fluoranthene	ND	0.0096	0.0050	mg/kg	
207-08-9	Benzo(k)fluoranthene	ND	0.0096	0.0050	mg/kg	
50-32-8	Benzo(a)pyrene	ND	0.0096	0.0050	mg/kg	
218-01-9	Chrysene	ND	0.0096	0.0050	mg/kg	
53-70-3	Dibenz(a,h)anthracene	ND	0.0096	0.0050	mg/kg	
206-44-0	Fluoranthene	ND	0.0096	0.0050	mg/kg	
86-73-7	Fluorene	ND	0.0096	0.0050	mg/kg	
193-39-5	Indeno(1,2,3-cd)pyrene	ND	0.0096	0.0050	mg/kg	
91-20-3	Naphthalene	0.0348	0.013	0.012	mg/kg	
129-00-0	Pyrene	ND	0.0096	0.0050	mg/kg	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
4165-60-0	Nitrobenzene-d5	80%		10-159%
321-60-8	2-Fluorobiphenyl	85%		19-131%
1718-51-0	Terphenyl-d14	85%		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

4

Accutest Laboratories

Report of Analysis

Page 1 of 1

Client Sample ID: CUT 2 SUBLINER (COMP)**Lab Sample ID:** D41014-1**Date Sampled:** 11/14/12**Matrix:** SO - Soil**Date Received:** 11/16/12**Method:** SW846 8015B**Percent Solids:** 86.4**Project:** PCU 296-5A

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	GB18507.D	1	11/16/12	SK	n/a	n/a	GGB1010
Run #2							

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

CAS No.	Compound	Result	RL	MDL	Units	Q
	TPH-GRO (C6-C10)	ND	13	6.5	mg/kg	
CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits		
120-82-1	1,2,4-Trichlorobenzene	92%		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: CUT 2 SUBLINER (COMP)**Lab Sample ID:** D41014-1**Date Sampled:** 11/14/12**Matrix:** SO - Soil**Date Received:** 11/16/12**Method:** SW846-8015B SW846 3546**Percent Solids:** 86.4**Project:** PCU 296-5A

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	FD19691.D	1	11/19/12	AV	11/19/12	OP6979	GFD990
Run #2							

	Initial Weight	Final Volume
Run #1	30.0 g	2.0 ml
Run #2		

CAS No.	Compound	Result	RL	MDL	Units	Q
	TPH-DRO (C10-C28)	24.7	15	10	mg/kg	
CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits		
84-15-1	o-Terphenyl	79%		35-130%		

ND = Not detected MDL - Method Detection Limit

J = Indicates an estimated value

RL = Reporting Limit

B = Indicates analyte found in associated method blank

E = Indicates value exceeds calibration range

N = Indicates presumptive evidence of a compound

4.1

4

Report of Analysis

Page 1 of 1

Client Sample ID: CUT 2 SUBLINER (COMP)**Lab Sample ID:** D41014-1**Matrix:** SO - Soil**Date Sampled:** 11/14/12**Date Received:** 11/16/12**Percent Solids:** 86.4**Project:** PCU 296-5A**Metals Analysis**

Analyte	Result	RL	Units	DF	Prep	Analyzed By	Method	Prep Method
Arsenic	13.0	0.11	mg/kg	5	11/19/12	11/26/12	JB	SW846 6020A ²
Barium	1710	1.1	mg/kg	1	11/19/12	11/19/12	JM	SW846 6010C ¹
Cadmium	< 1.1	1.1	mg/kg	1	11/19/12	11/19/12	JM	SW846 6010C ¹
Chromium	65.4	1.1	mg/kg	1	11/19/12	11/19/12	JM	SW846 6010C ¹
Copper	8.1	1.1	mg/kg	1	11/19/12	11/19/12	JM	SW846 6010C ¹
Lead	9.1	5.7	mg/kg	1	11/19/12	11/20/12	JM	SW846 6010C ¹
Mercury	< 0.090	0.090	mg/kg	1	11/27/12	11/27/12	JM	SW846 7471B ³
Nickel	20.5	3.4	mg/kg	1	11/19/12	11/19/12	JM	SW846 6010C ¹
Selenium	< 5.7	5.7	mg/kg	1	11/19/12	11/19/12	JM	SW846 6010C ¹
Silver	< 3.4	3.4	mg/kg	1	11/19/12	11/19/12	JM	SW846 6010C ¹
Zinc	38.2	3.4	mg/kg	1	11/19/12	11/19/12	JM	SW846 6010C ¹

(1) Instrument QC Batch: MA3012

(2) Instrument QC Batch: MA3019

(3) Instrument QC Batch: MA3023

(4) Prep QC Batch: MP8913

(5) Prep QC Batch: MP8914

(6) Prep QC Batch: MP8936

RL = Reporting Limit

Report of Analysis

Page 1 of 1

Client Sample ID: CUT 2 SUBLINER (COMP)**Lab Sample ID:** D41014-1**Matrix:** SO - Soil**Date Sampled:** 11/14/12**Date Received:** 11/16/12**Percent Solids:** 86.4**Project:** PCU 296-5A**General Chemistry**

Analyte	Result	RL	Units	DF	Analyzed	By	Method
prep: DEPT.OF AG, BOOK N9							
Specific Conductivity	353	1.0	umhos/cm	1	11/19/12	JD	SM2510B-1997 MOD
Chromium, Hexavalent	< 1.0	1.0	mg/kg	1	11/19/12	KB	SW846 3060A/7196A
Chromium, Trivalent ^a	65.4	2.1	mg/kg	1	11/19/12 17:50	JM	SW846 3060/7196A M
Redox Potential Vs H2	94.2		mv	1	11/16/12	CT	ASTM D1498-76M
Solids, Percent	86.4		%	1	11/19/12	SWT	SM19 2540B M
pH	9.90		su	1	11/16/12 15:45	JD	SW846 9045D

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

RL = Reporting Limit

Report of Analysis

Page 1 of 1

4.2
4**Client Sample ID:** CUT 2 SUBLINER (COMP)**Lab Sample ID:** D41014-1A**Matrix:** SO - Soil**Project:** PCU 296-5A**Date Sampled:** 11/14/12**Date Received:** 11/16/12**Percent Solids:** 86.4**SAR Metals Analysis**

Analyte	Result	RL	Units	DF	Prep	Analyzed By	Method	Prep Method
Calcium	14.0	2.0	mg/l	1	11/19/12	11/19/12 JM	SW846 6010C ¹	SW846 3010A/M ²
Magnesium	4.53	1.0	mg/l	1	11/19/12	11/19/12 JM	SW846 6010C ¹	SW846 3010A/M ²
Sodium	61.8	2.0	mg/l	1	11/19/12	11/19/12 JM	SW846 6010C ¹	SW846 3010A/M ²

(1) Instrument QC Batch: MA3012

(2) Prep QC Batch: MP8915

RL = Reporting Limit

Report of Analysis

Page 1 of 1

4.2
4**Client Sample ID:** CUT 2 SUBLINER (COMP)**Lab Sample ID:** D41014-1A**Matrix:** SO - Soil**Project:** PCU 296-5A**Date Sampled:** 11/14/12**Date Received:** 11/16/12**Percent Solids:** 86.4**General Chemistry**

Analyte	Result	RL	Units	DF	Analyzed	By	Method
Sodium Adsorption Ratio ^a	3.67		ratio	1	11/19/12 15:32	JM	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.accurtest.com

Client / Reporting Information		Project Information		Requested Analysis (see TEST CODE sheet)										Matrix Codes																													
Company Name KRW Consulting	Project Name: XTO PCU 296-5A																																										
Street Address 8000 West 14th Street; Suite 200	Street																																										
City Lakewood, CO 80214	City	Billing Information (if different from Report to)																																									
Project Contact Dwayne Knudson	Project # 1210-04	Company Name XTO Energy																																									
Phone # 970-488-1098	Client Purchase Order #	Street Address 21459 CR 5																																									
Sampler(s) Name(s) DAVID SANDERS	Project Manager Joe Hess	City Rifle, CO 81650																																									
Sampler(s) Name(s) DAVID SANDERS	Project Manager Joe Hess	Attention: Jessica Dooling																																									
Accutest Sample #		Field ID / Point of Collection CUT 2SUBLINER (COMP)		Collection			Sampled by	Number of preserved Bottles						T-910	LAB USE ONLY O1																												
				Date 11-14-12	Time 13:45	METH/OD Vial # DS		# of bottles 50	High 1	Med 0	Low 0	None 0	Unknown 0			MEOH 0	ENCORE 0	Bisulfite 0																									
Turnaround Time (Business days)		Data Deliverable Information										Comments / Special Instructions																															
<input type="checkbox"/> Std. 10 Business Days <input checked="" type="checkbox"/> Std. 5 Business Days (By contract only) <input type="checkbox"/> 3 Day Emergency <input type="checkbox"/> 2 Day Emergency <input type="checkbox"/> 1 Day Emergency <input type="checkbox"/>		Approved By (Accutest PM): Date: _____ _____ _____ _____ _____										<input type="checkbox"/> Commercial "A" (Level 1) <input type="checkbox"/> Commercial "B" (Level 2) <input type="checkbox"/> COMM BN <input type="checkbox"/> COMM BN+ <input type="checkbox"/>			<input type="checkbox"/> State Forms Required <input type="checkbox"/> Send Forms to State <input type="checkbox"/> Report by Fax <input checked="" type="checkbox"/> Report by PDF ONLY <input type="checkbox"/> EDD Format																												
Emergency & Rush T/A data available VIA Lablink												Please email to: KRW Piceance Team																															
<p>Sample Custody must be documented below each time samples change possession, including courier delivery.</p> <table border="1"> <tr> <td>Relinquished by Sampler: 1 Lori Abu-Saad</td> <td>Date Time: 11/15/12 17:00</td> <td>Received By: Rifle Service Center</td> <td>Relinquished By: 2</td> <td>Date Time:</td> <td>Received By:</td> </tr> <tr> <td>Relinquished by Sampler: 3</td> <td>Date Time:</td> <td>Received By: 3</td> <td>Relinquished By: 4</td> <td>Date Time:</td> <td>Received By: 2</td> </tr> <tr> <td>Relinquished by: 5</td> <td>Date Time:</td> <td>Received By: 5</td> <td>Custody Seal # HPCO</td> <td>Intact <input type="checkbox"/></td> <td>Preserved where applicable <input type="checkbox"/></td> </tr> <tr> <td></td> <td></td> <td></td> <td></td> <td>Not Intact <input type="checkbox"/></td> <td>Orals <input type="checkbox"/></td> </tr> <tr> <td></td> <td></td> <td></td> <td></td> <td></td> <td>Cooler Temp. 27.1 C</td> </tr> </table>														Relinquished by Sampler: 1 Lori Abu-Saad	Date Time: 11/15/12 17:00	Received By: Rifle Service Center	Relinquished By: 2	Date Time:	Received By:	Relinquished by Sampler: 3	Date Time:	Received By: 3	Relinquished By: 4	Date Time:	Received By: 2	Relinquished by: 5	Date Time:	Received By: 5	Custody Seal # HPCO	Intact <input type="checkbox"/>	Preserved where applicable <input type="checkbox"/>					Not Intact <input type="checkbox"/>	Orals <input type="checkbox"/>						Cooler Temp. 27.1 C
Relinquished by Sampler: 1 Lori Abu-Saad	Date Time: 11/15/12 17:00	Received By: Rifle Service Center	Relinquished By: 2	Date Time:	Received By:																																						
Relinquished by Sampler: 3	Date Time:	Received By: 3	Relinquished By: 4	Date Time:	Received By: 2																																						
Relinquished by: 5	Date Time:	Received By: 5	Custody Seal # HPCO	Intact <input type="checkbox"/>	Preserved where applicable <input type="checkbox"/>																																						
				Not Intact <input type="checkbox"/>	Orals <input type="checkbox"/>																																						
					Cooler Temp. 27.1 C																																						

5.1

D41014: Chain of Custody



Accutest Laboratories Sample Receipt Summary

Accutest Job Number: D41014

Client: KRW CONSULTING

Immediate Client Services Action Required: No

Date / Time Received: 11/16/2012 1:00:00 PM

No. Coolers:

1

Client Service Action Required at Login: No

Project: XTO PCU 296-5A

Airbill #'s: HDCO

Cooler Security Y or N

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

Cooler Temperature Y or N

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

Quality Control Preservation Y or N N/A

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

Sample Integrity - Documentation

- | | | |
|--|-------------------------------------|--------------------------|
| 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 rcvd 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

D41014: Chain of Custody

Page 2 of 2



GC/MS Volatiles

QC Data Summaries

Includes the following where applicable:

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



Method Blank Summary

Job Number: D41014

Account: XTOKWR XTO Energy

Project: PCU 296-5A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
V5V1506-MB	5V24673.D	1	11/19/12	BD	n/a	n/a	V5V1506

The QC reported here applies to the following samples:

Method: SW846 8260B

D41014-1

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	101%
460-00-4	4-Bromofluorobenzene	93%
17060-07-0	1,2-Dichloroethane-D4	95%

Blank Spike Summary

Job Number: D41014
Account: XTOKWR XTO Energy
Project: PCU 296-5A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
V5V1506-BS	5V24674.D	1	11/19/12	BD	n/a	n/a	V5V1506

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

D41014-1

CAS No.	Compound	Spike ug/kg	BSP ug/kg	BSP %	Limits
71-43-2	Benzene	50	54.0	108	70-130
100-41-4	Ethylbenzene	50	54.2	108	70-130
108-88-3	Toluene	50	53.2	106	70-130
1330-20-7	Xylene (total)	150	168	112	70-130

CAS No.	Surrogate Recoveries	BSP	Limits
2037-26-5	Toluene-D8	100%	64-130%
460-00-4	4-Bromofluorobenzene	98%	62-131%
17060-07-0	1,2-Dichloroethane-D4	96%	70-130%

* = Outside of Control Limits.

Matrix Spike/Matrix Spike Duplicate Summary

Page 1 of 1

Job Number: D41014

Account: XTOKWR XTO Energy

Project: PCU 296-5A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
D41036-1MS	5V24676.D	1	11/19/12	BD	n/a	n/a	V5V1506
D41036-1MSD	5V24677.D	1	11/19/12	BD	n/a	n/a	V5V1506
D41036-1	5V24675.D	1	11/19/12	BD	n/a	n/a	V5V1506

The QC reported here applies to the following samples:

Method: SW846 8260B

D41014-1

CAS No.	Compound	D41036-1		Spike	MS	MS	MSD	MSD	Limits	
		ug/kg	Q	ug/kg	ug/kg	%	ug/kg	%	RPD	Rec/RPD
71-43-2	Benzene	702		5720	6570	103	6760	106	3	64-139/30
100-41-4	Ethylbenzene	2750		5720	8410	99	8720	104	4	68-136/30
108-88-3	Toluene	2340		5720	7850	96	7970	98	2	60-130/30
1330-20-7	Xylene (total)	21000		17200	35900	87	37500	96	4	58-142/30

CAS No.	Surrogate Recoveries	MS	MSD	D41036-1	Limits
2037-26-5	Toluene-D8	99%	98%	100%	64-130%
460-00-4	4-Bromofluorobenzene	105%	108%	103%	62-131%
17060-07-0	1,2-Dichloroethane-D4	97%	98%	95%	70-130%

* = Outside of Control Limits.

6.3.1
6



GC/MS Volatiles

Raw Data

7

Judy Nelson
 11/20/12 12:37

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\V5111912.S\
 Data File : 5V24680.D
 Acq On : 19 Nov 2012 4:33 pm
 Operator : BRETD
 Sample : D41014-1
 Misc : MS4990,V5V1506,5.068,,100,5,1
 ALS Vial : 12 Sample Multiplier: 1

Quant Time: Nov 20 08:58:55 2012
 Quant Method : C:\msdchem\1\METHODS\V5AP1497TVH1497.M
 Quant Title : 8260
 QLast Update : Wed Nov 14 09:54:38 2012
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
2) Pentafluorobenzene	11.624	168	442147	50.00	ug/l	0.00
35) 1,4-Difluorobenzene	12.423	114	536703	50.00	ug/l	0.00
53) Chlorobenzene-d5	15.072	117	501261	50.00	ug/l	0.00
74) 1,4-Dichlorobenzene-d4	17.036	152	375149	50.00	ug/l	0.00

System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	12.024	102	36310	48.25	ug/l	0.00
Spiked Amount 50.000	Range 70 - 130		Recovery	=	96.50%	
61) Toluene-d8	13.816	98	597215	50.29	ug/l	0.00
Spiked Amount 50.000	Range 70 - 130		Recovery	=	100.58%	
69) 4-Bromofluorobenzene	16.020	95	255528	49.93	ug/l	0.00
Spiked Amount 50.000	Range 70 - 130		Recovery	=	99.86%	

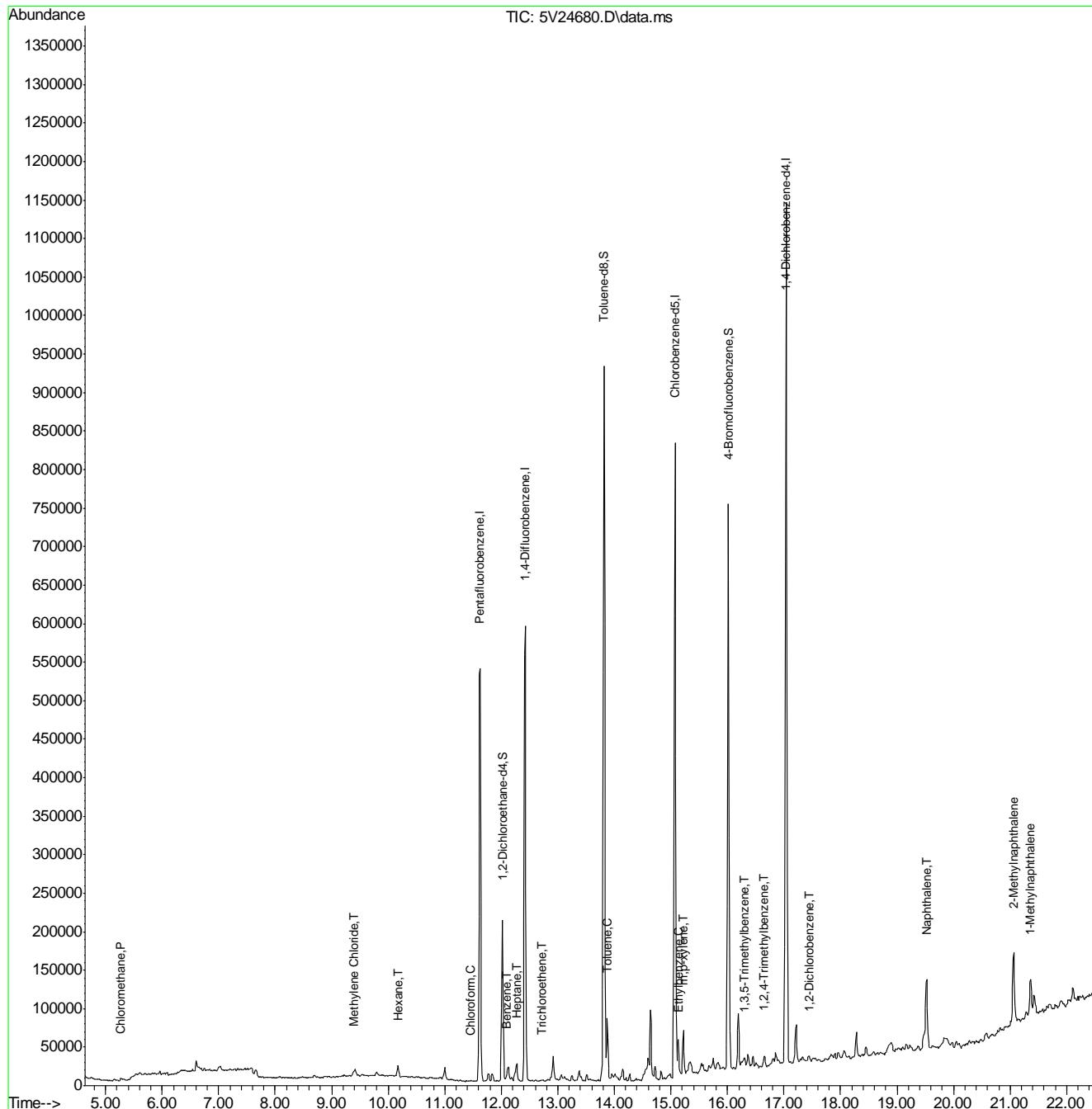
Target Compounds					Qvalue
4) Chloromethane	5.265	50	1771	0.36	ug/l 100
17) Methylene Chloride	9.387	84	2041	0.51	ug/l # 80
29) Chloroform	11.453	83	384	0.06	ug/l 75
41) Hexane	10.174	57	6721	1.38	ug/l 100
43) Heptane	12.275	43	9325	1.67	ug/l 92
48) Trichloroethene	12.709	95	427	0.11	ug/l # 86
50) Benzene	12.092	78	9734	0.68	ug/l 100
62) Toluene	13.873	92	22472	2.43	ug/l 98
66) Ethylbenzene	15.141	91	8987	0.51	ug/l 91
72) m,p-xylene	15.221	106	17579	2.48	ug/l 96
80) 1,3,5-Trimethylbenzene	16.305	105	4098m	0.24	ug/l
82) 1,2,4-Trimethylbenzene	16.648	105	7525	0.42	ug/l 90
87) 1,2-Dichlorobenzene	17.447	146	462	0.04	ug/l # 78
91) Naphthalene	19.525	128	115642	6.22	ug/l 100
94) 2-Methylnaphthalene	21.066	142	67966	10.95	ug/l 97
95) 1-Methylnaphthalene	21.363	142	37759	4.85	ug/l 95

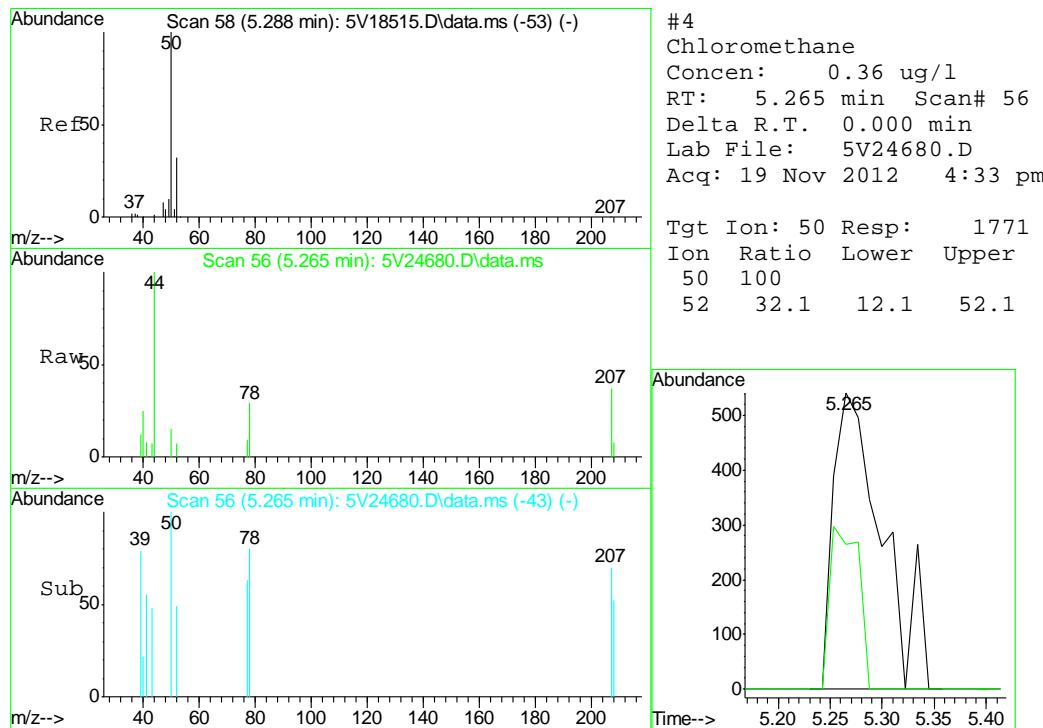
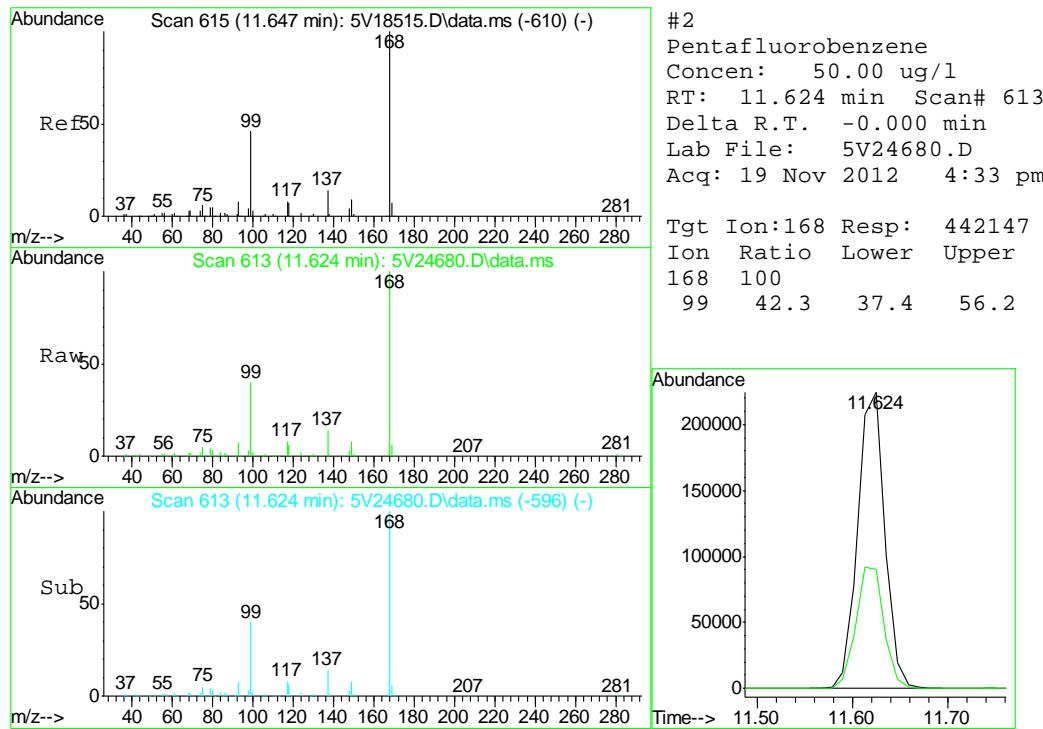
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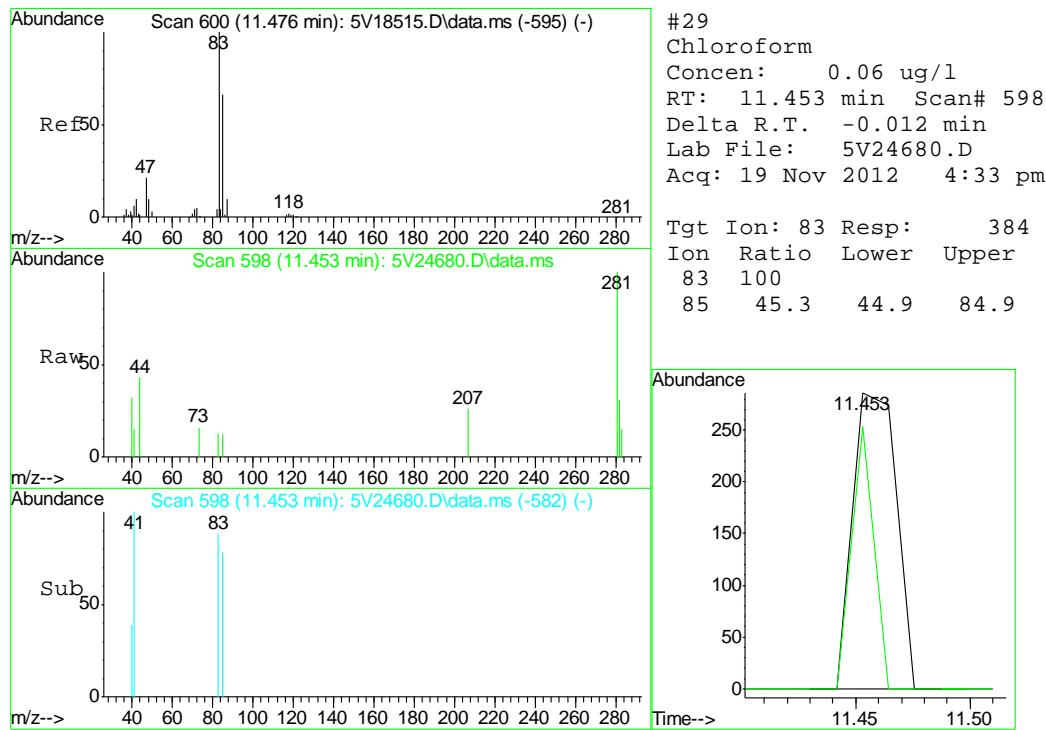
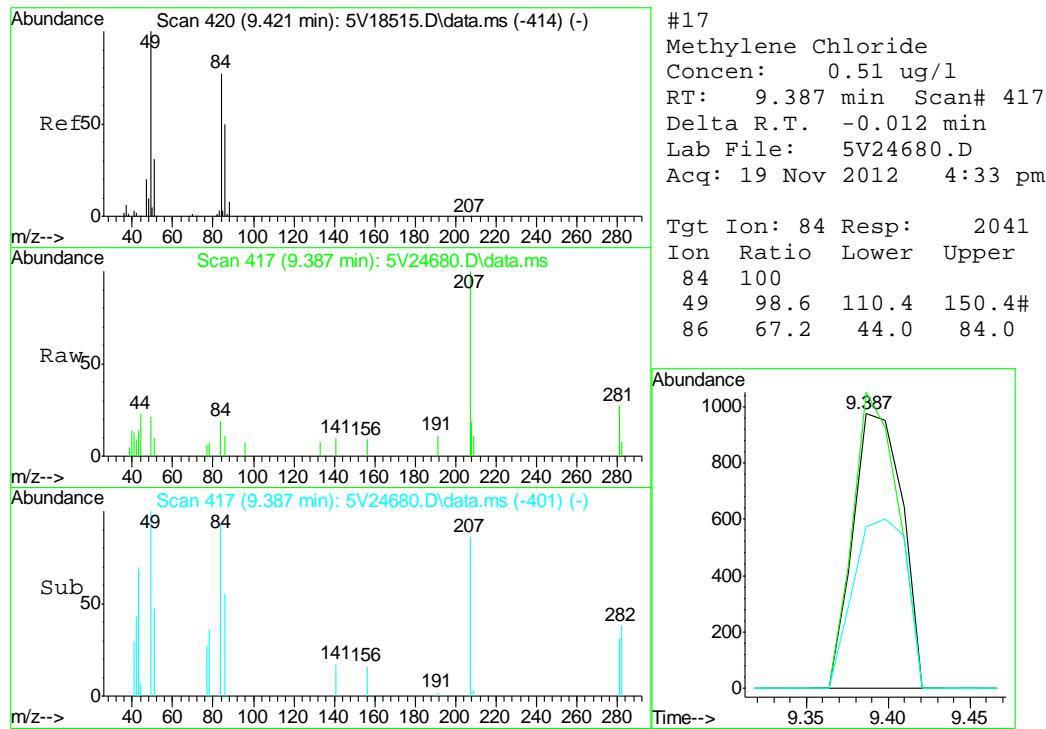
Quantitation Report (QT Reviewed)

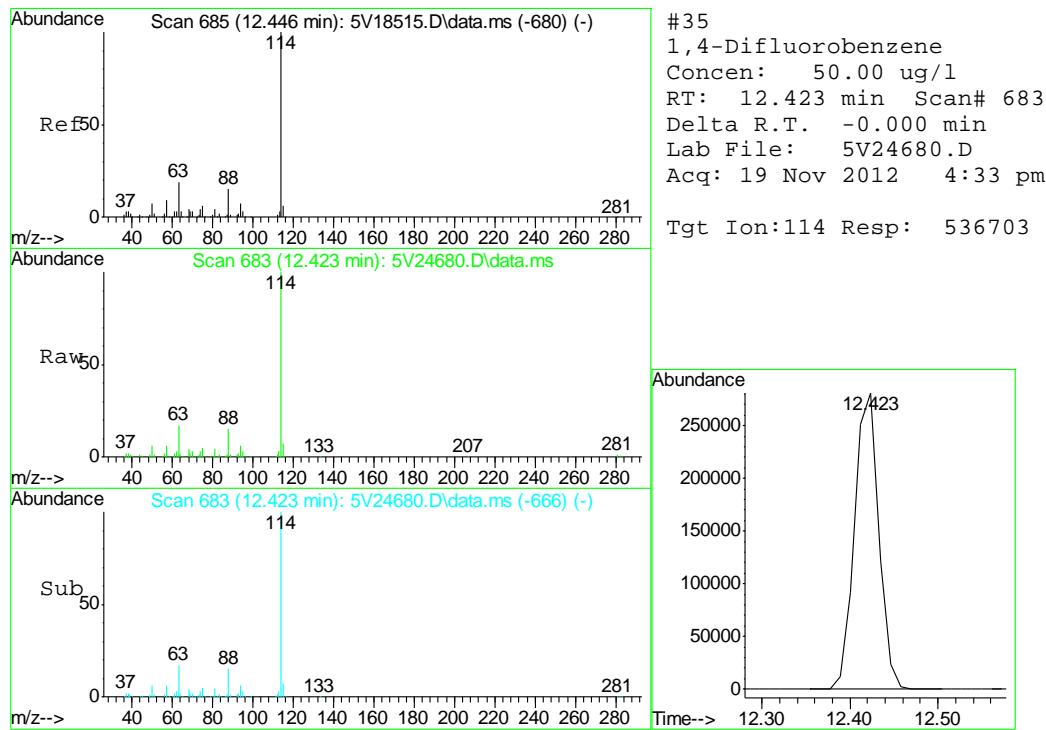
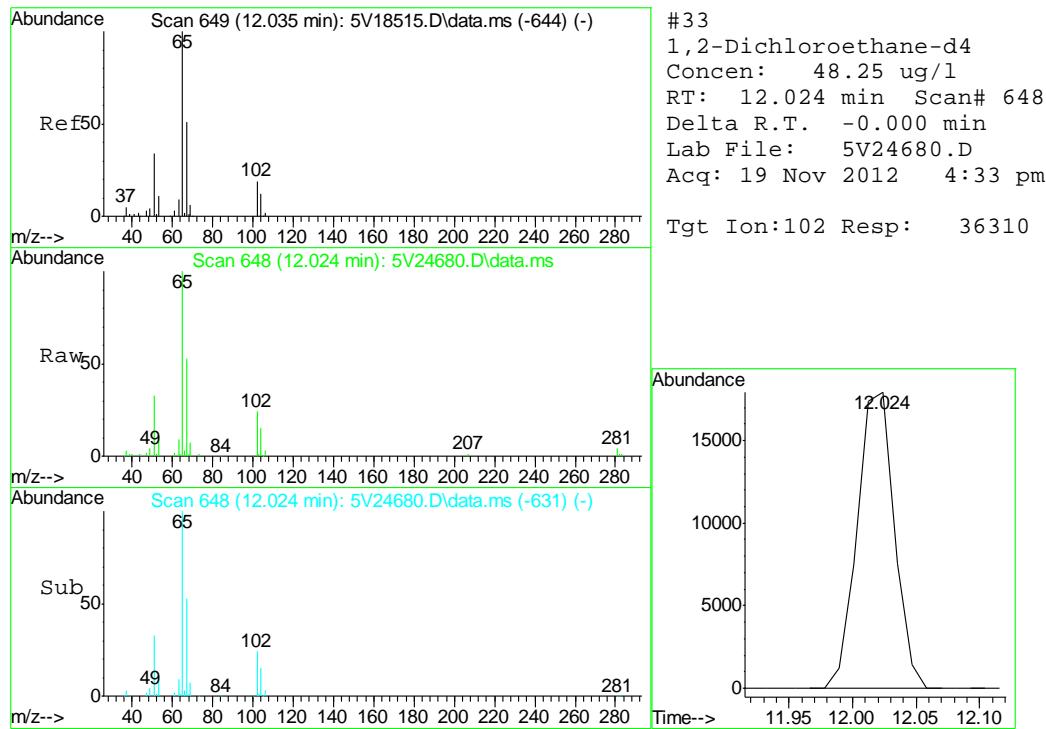
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 Data File : 5V24680.D
 Acq On : 19 Nov 2012 4:33 pm
 Operator : BRETD
 Sample : D41014-1
 Misc : MS4990,V5V1506,,5.068,,100,,5,1
 ALS Vial : 12 Sample Multiplier: 1

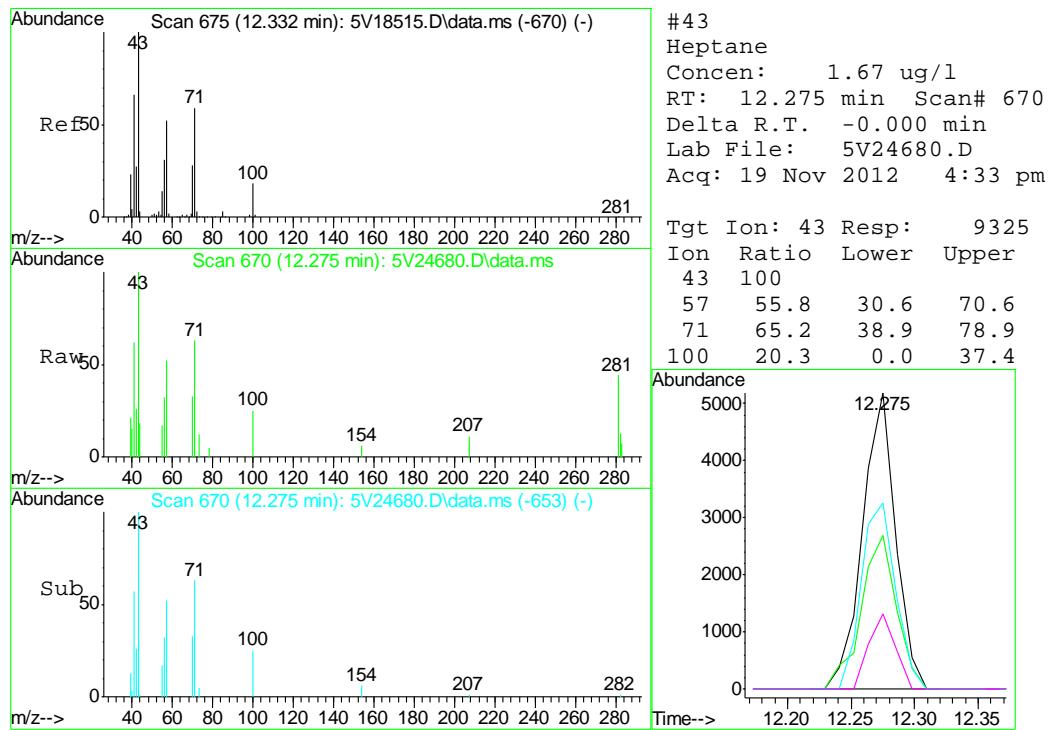
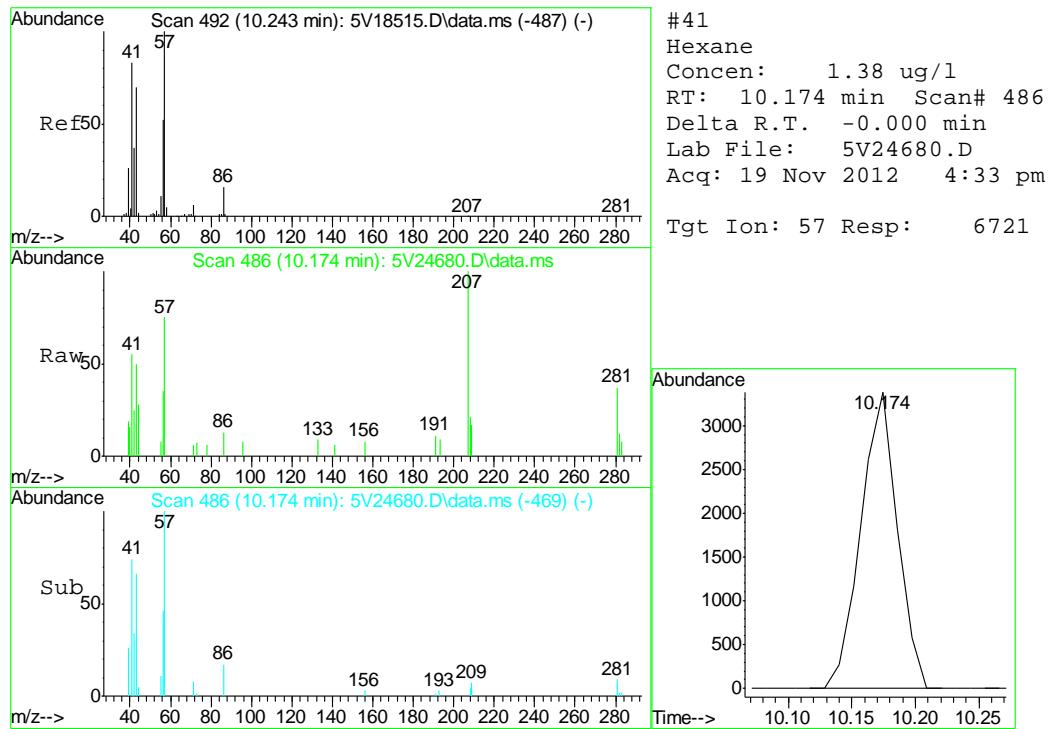
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 Quant Title : 8260
 QLast Update : Wed Nov 14 09:54:38 2012
 Response via : Initial Calibration

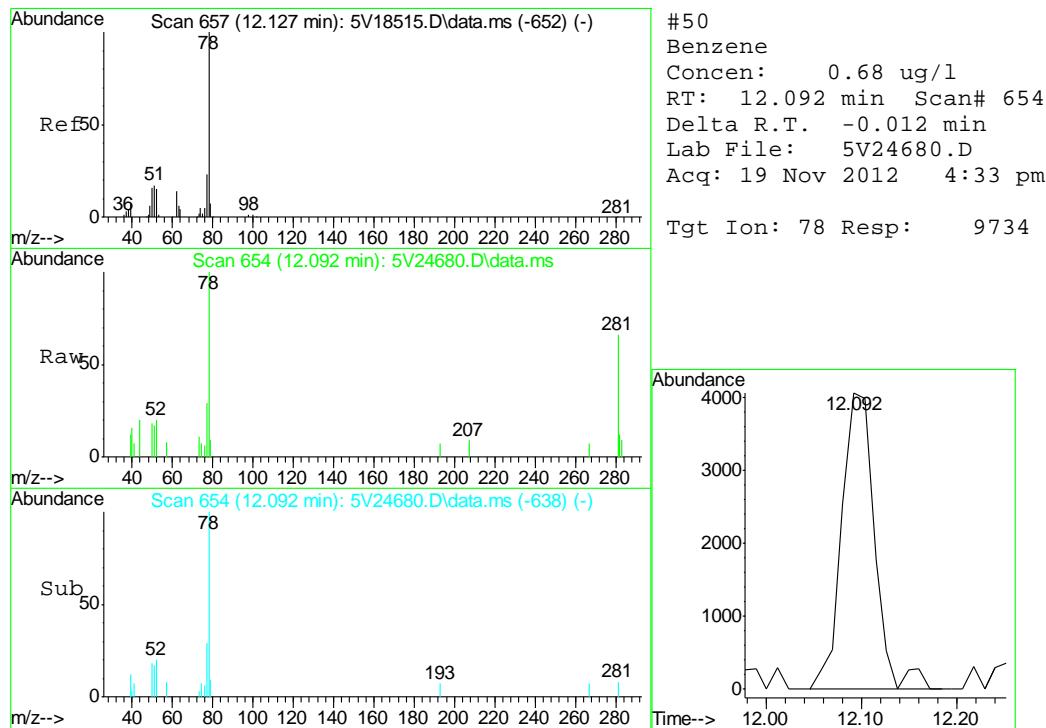
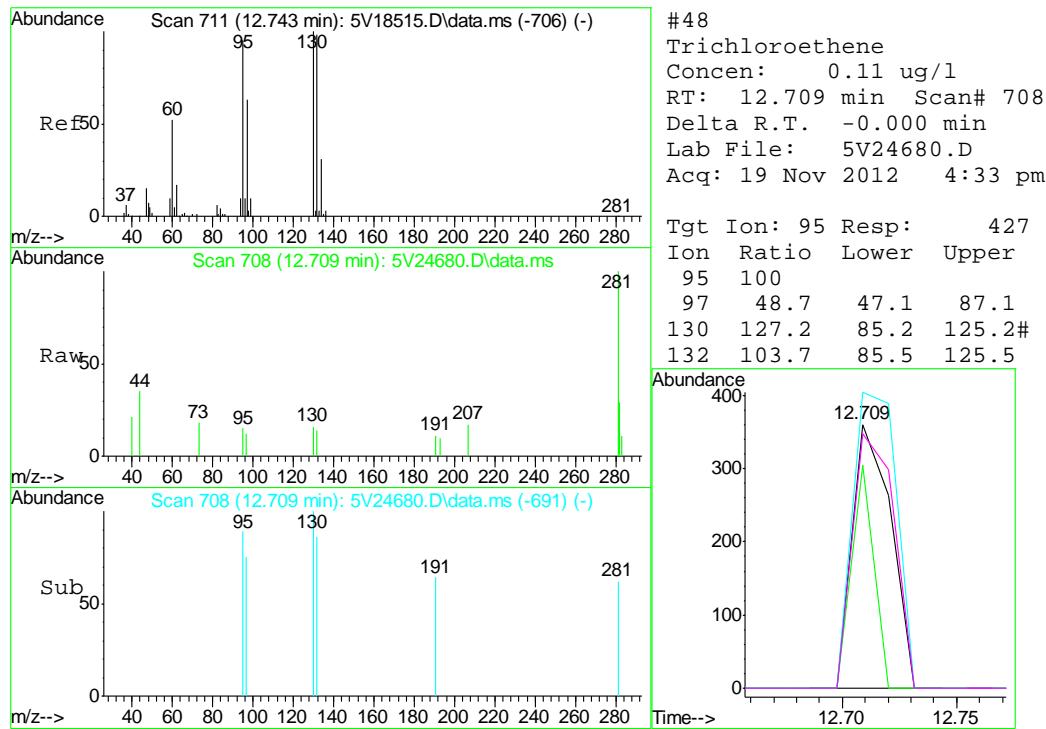


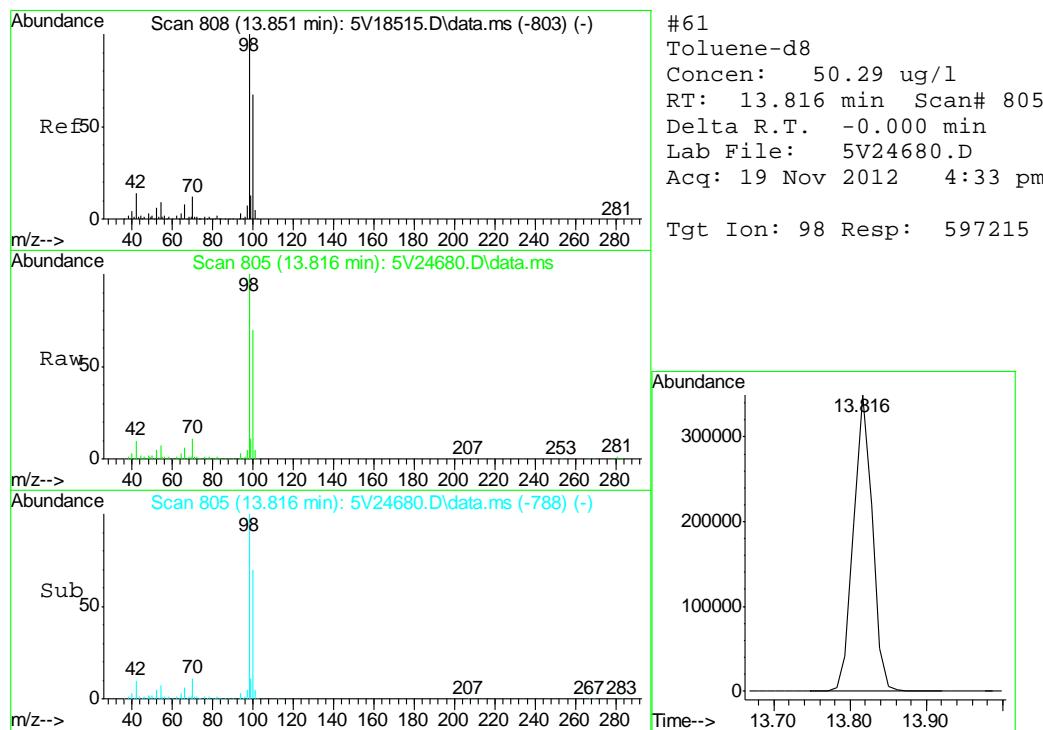
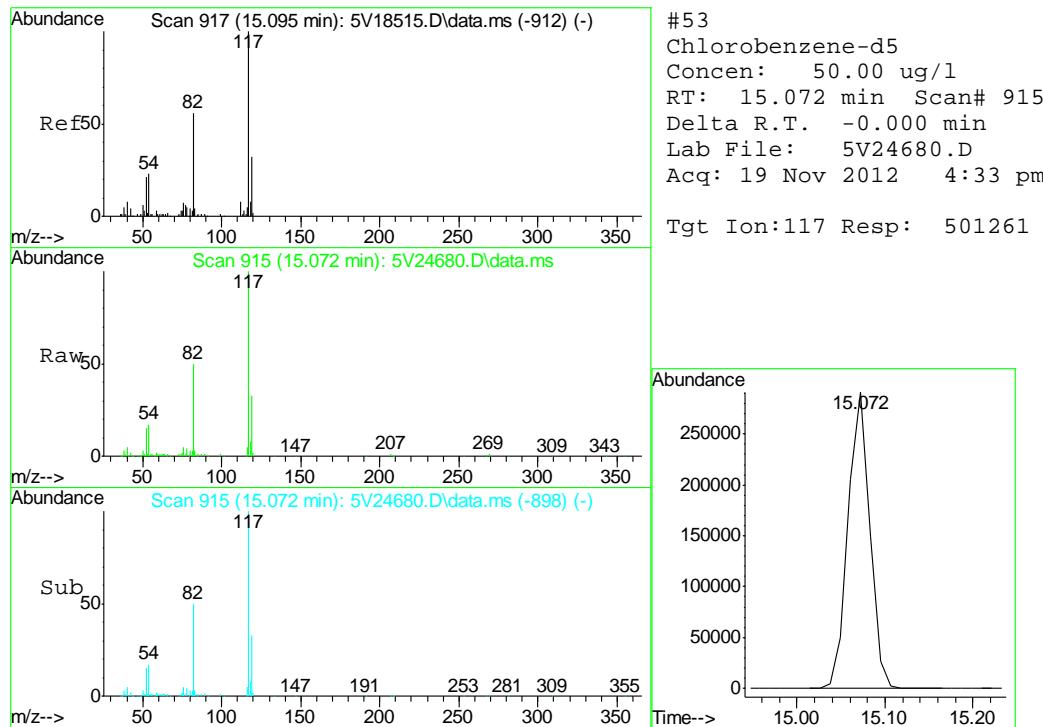


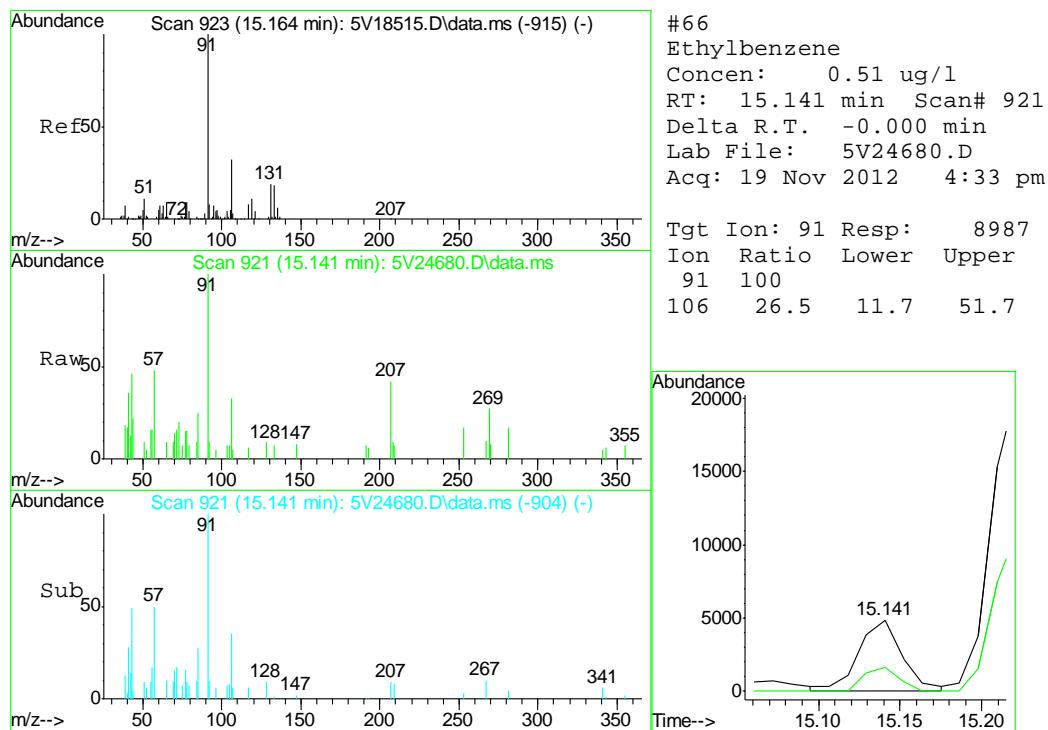
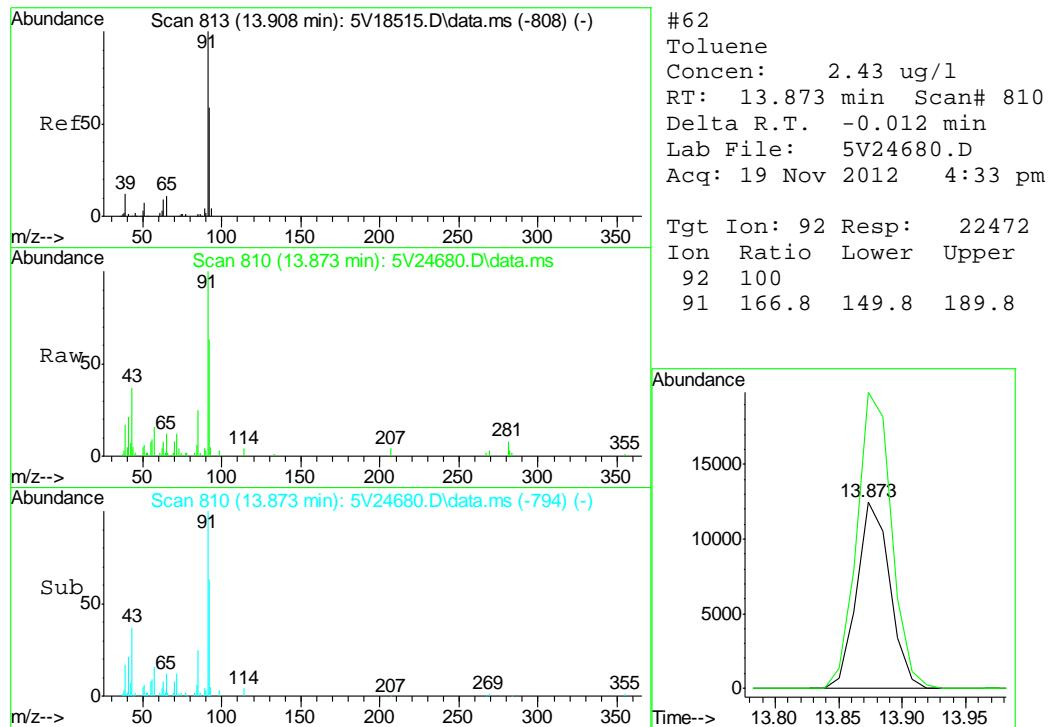


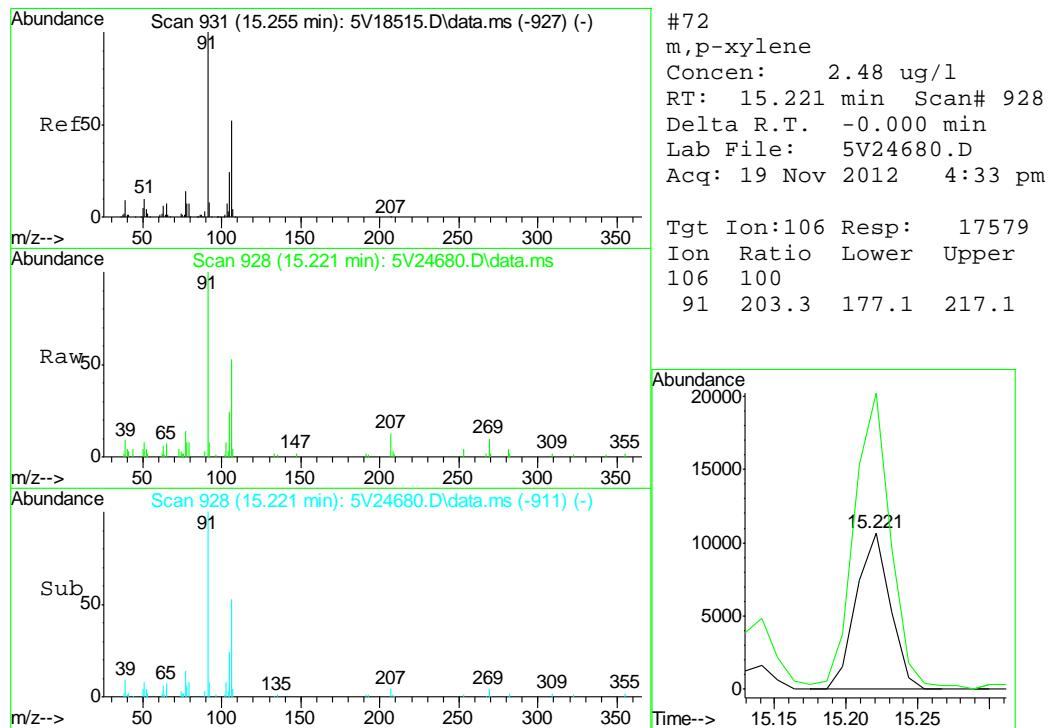
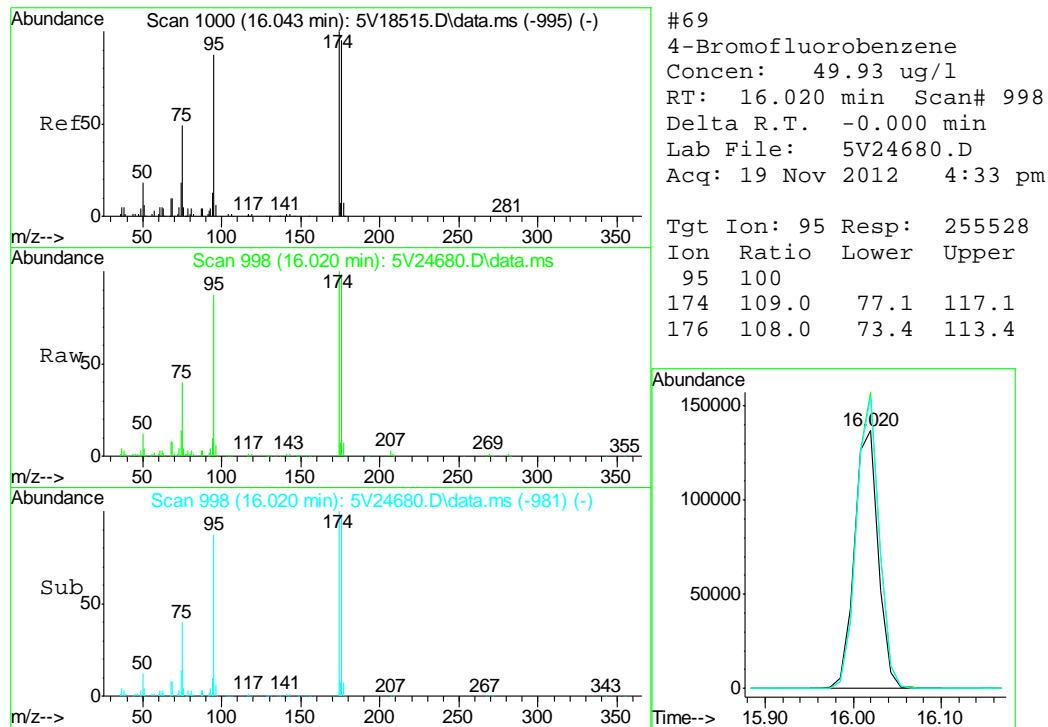


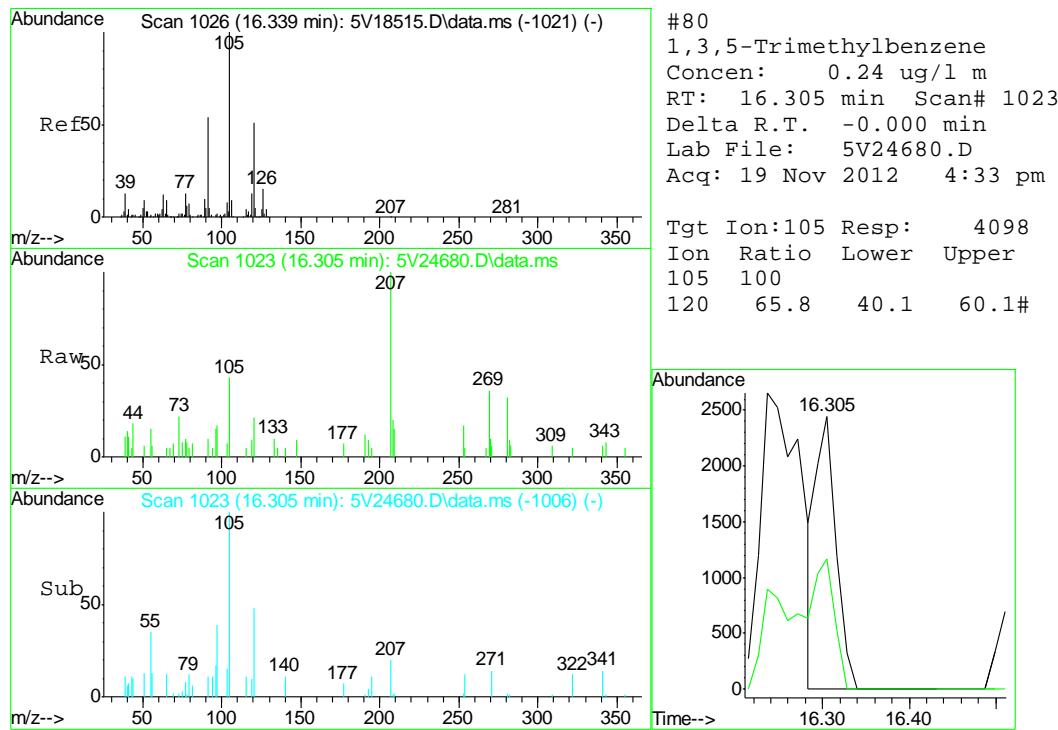
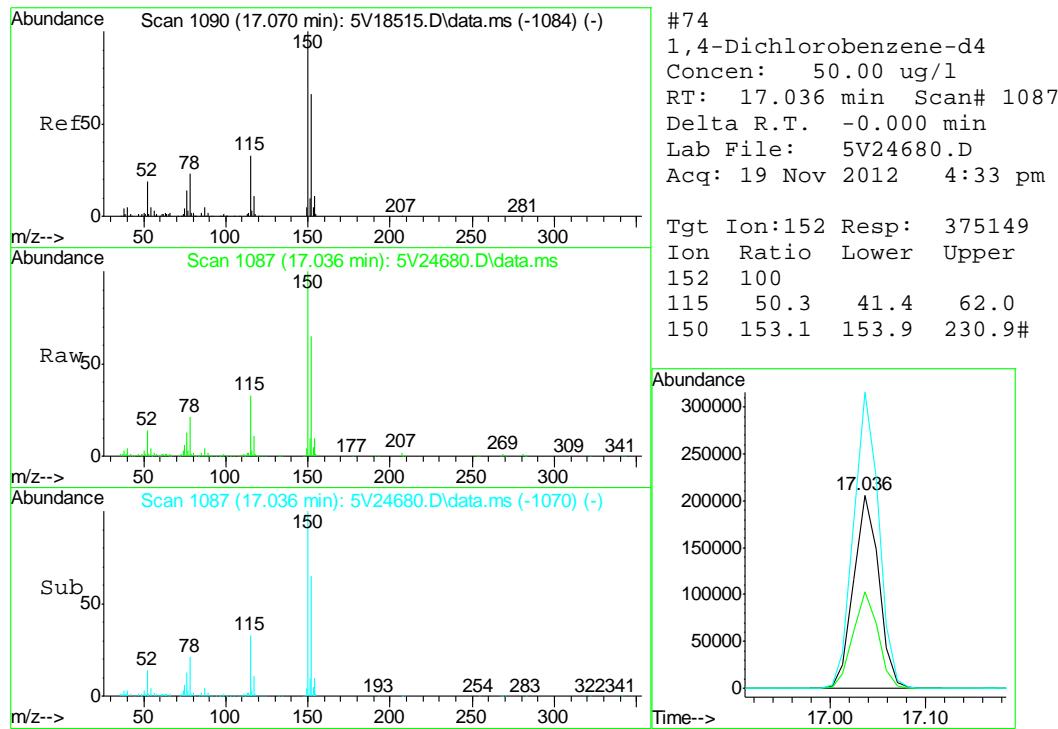


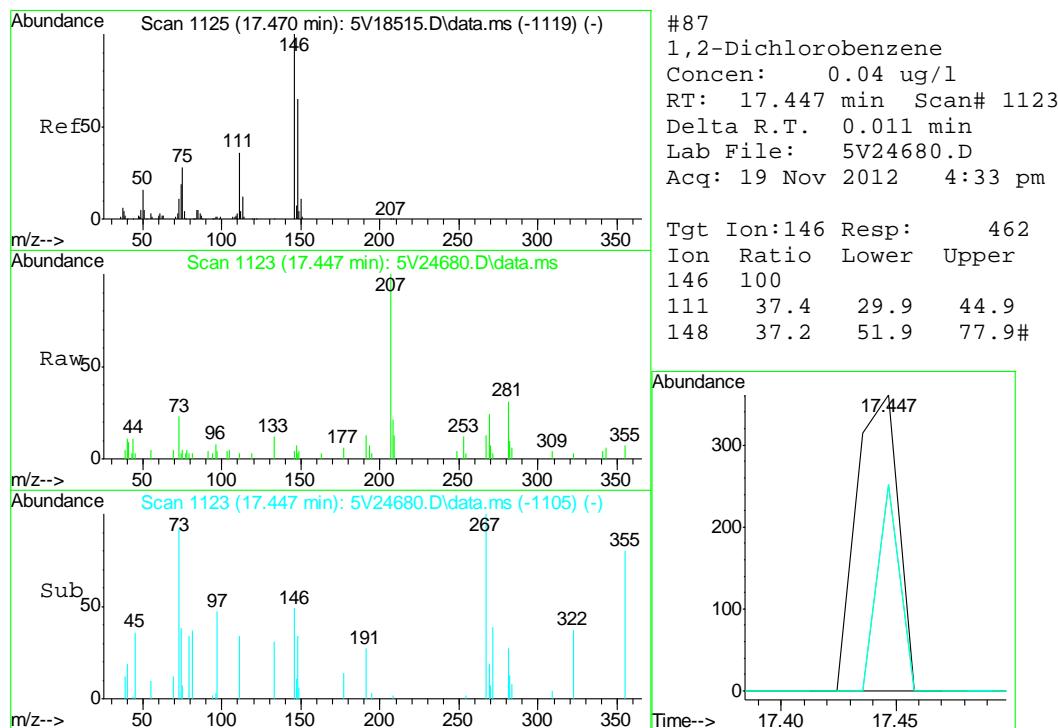
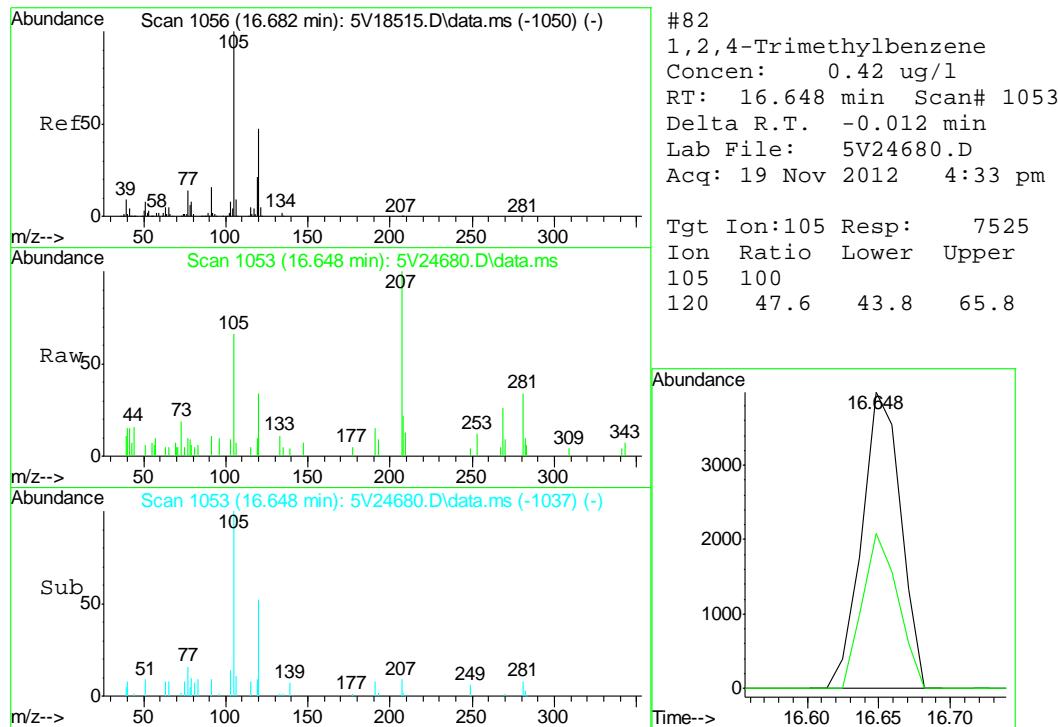


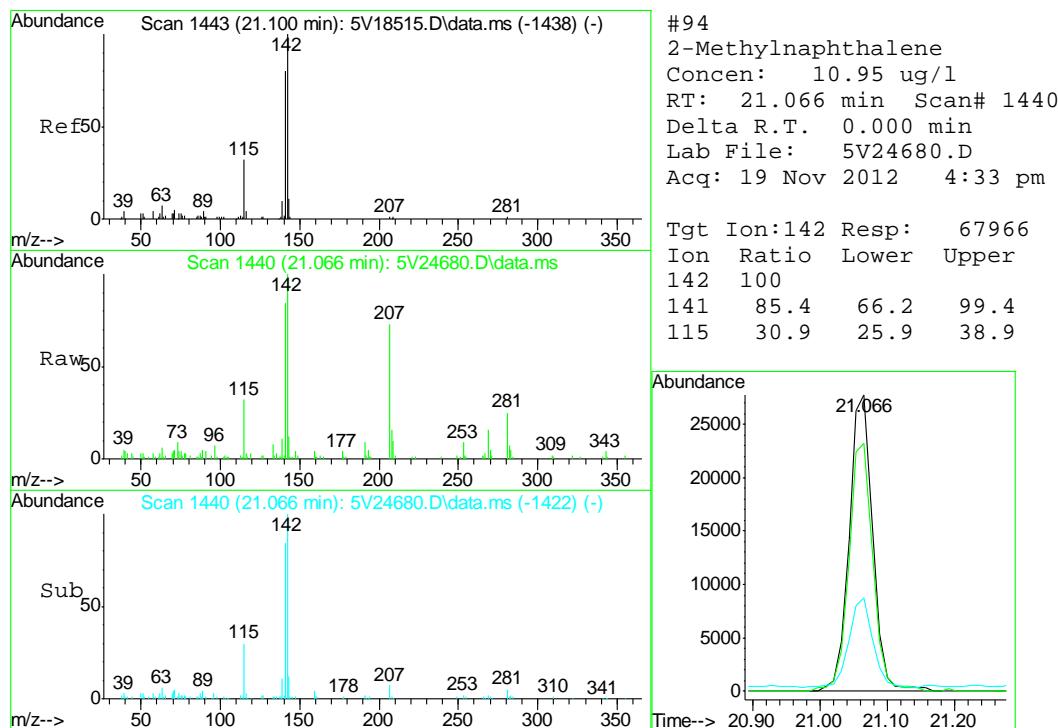
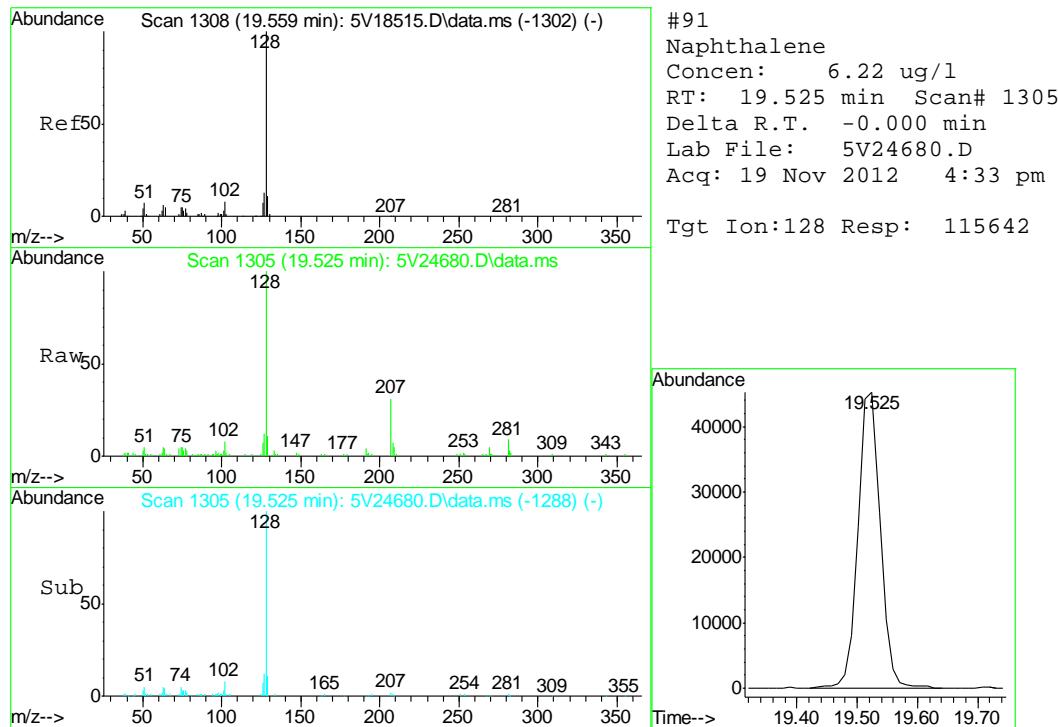


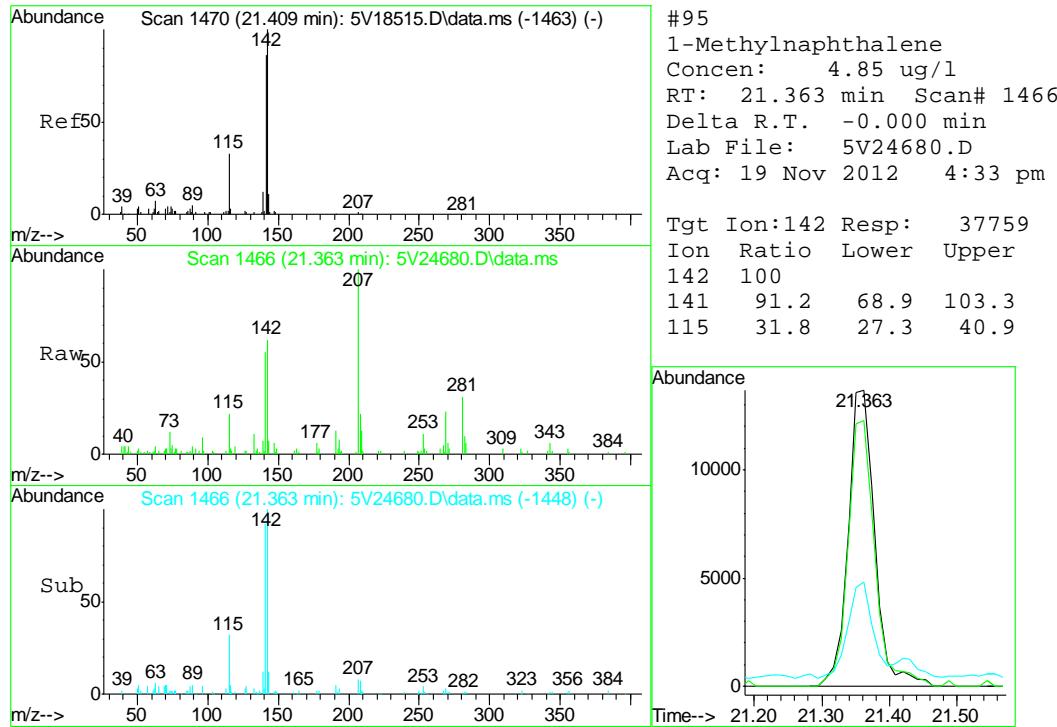












Quantitation Report (QT Reviewed)

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 Operator : BRETD
 Sample : MB
 Misc : MS4990,V5V1506,5.00,,100,5,1
 ALS Vial : 5 Sample Multiplier: 1

Quant Time: Nov 20 08:43:29 2012
 Quant Method : C:\msdchem\1\METHODS\V5AP1497TVH1497.M
 Quant Title : 8260
 QLast Update : Wed Nov 14 09:54:38 2012
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
2) Pentafluorobenzene	11.624	168	473270	50.00	ug/l	0.00
35) 1,4-Difluorobenzene	12.423	114	576356	50.00	ug/l	0.00
53) Chlorobenzene-d5	15.072	117	535944	50.00	ug/l	0.00
74) 1,4-Dichlorobenzene-d4	17.036	152	360659	50.00	ug/l	0.00

System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	12.024	102	38220	47.45	ug/l	0.00
Spiked Amount	50.000	Range	70 - 130	Recovery	=	94.90%
61) Toluene-d8	13.816	98	640357	50.44	ug/l	0.00
Spiked Amount	50.000	Range	70 - 130	Recovery	=	100.88%
69) 4-Bromofluorobenzene	16.020	95	254413	46.49	ug/l	0.00
Spiked Amount	50.000	Range	70 - 130	Recovery	=	92.98%

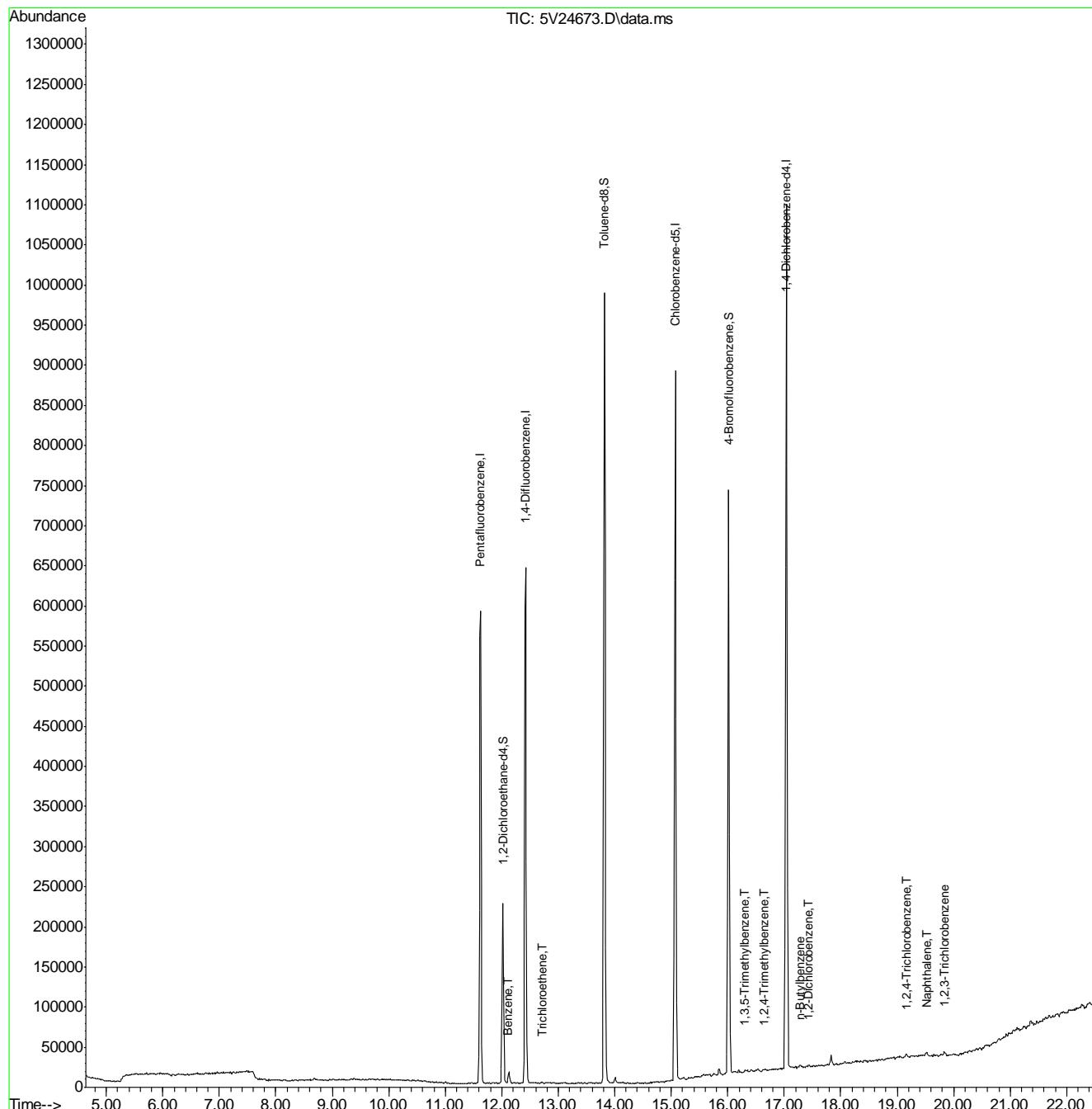
Target Compounds					Qvalue
48) Trichloroethene	12.709	95	778	0.19	ug/l
50) Benzene	12.104	78	1232	0.08	ug/l
80) 1,3,5-Trimethylbenzene	16.305	105	800	0.05	ug/l
82) 1,2,4-Trimethylbenzene	16.648	105	1132	0.07	ug/l
87) 1,2-Dichlorobenzene	17.436	146	1204	0.11	ug/l
88) n-Butylbenzene	17.287	91	2178	0.12	ug/l
90) 1,2,4-Trichlorobenzene	19.159	180	2308	0.23	ug/l
91) Naphthalene	19.525	128	6488	0.36	ug/l
93) 1,2,3-Trichlorobenzene	19.833	180	2556	0.27	ug/l

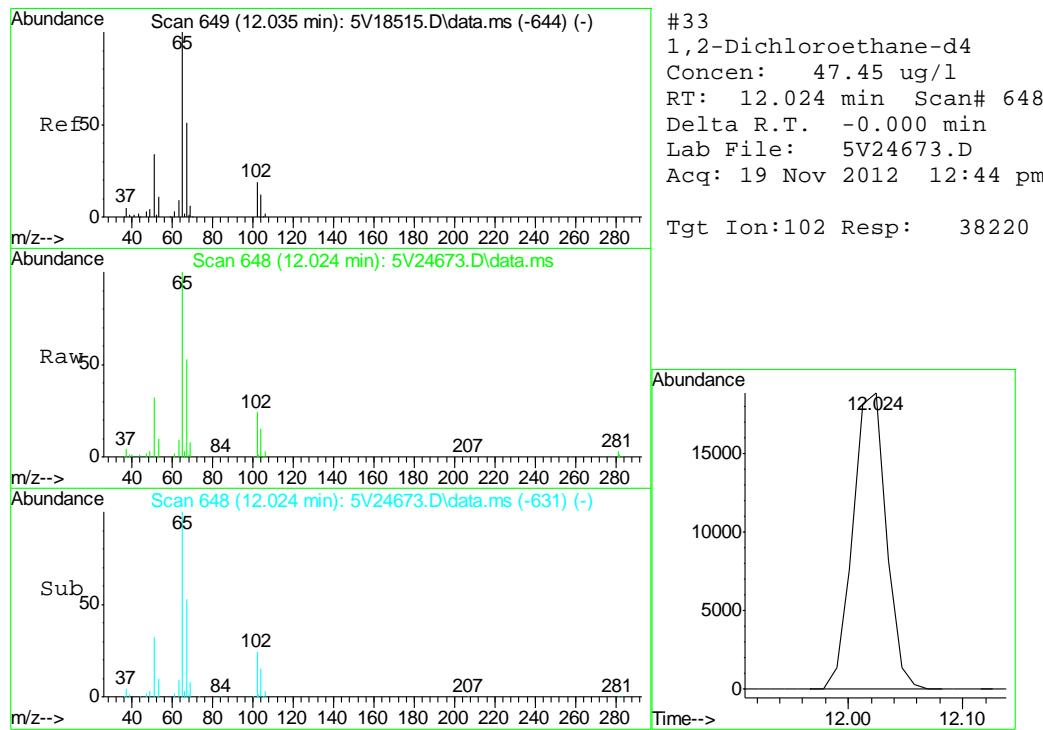
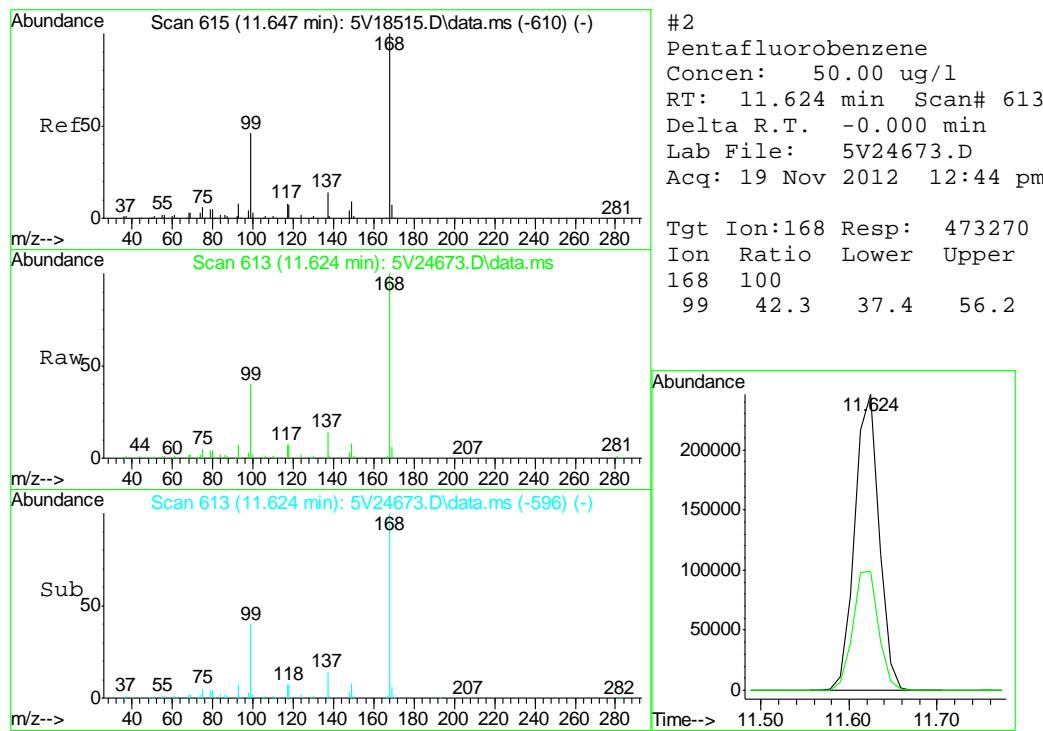
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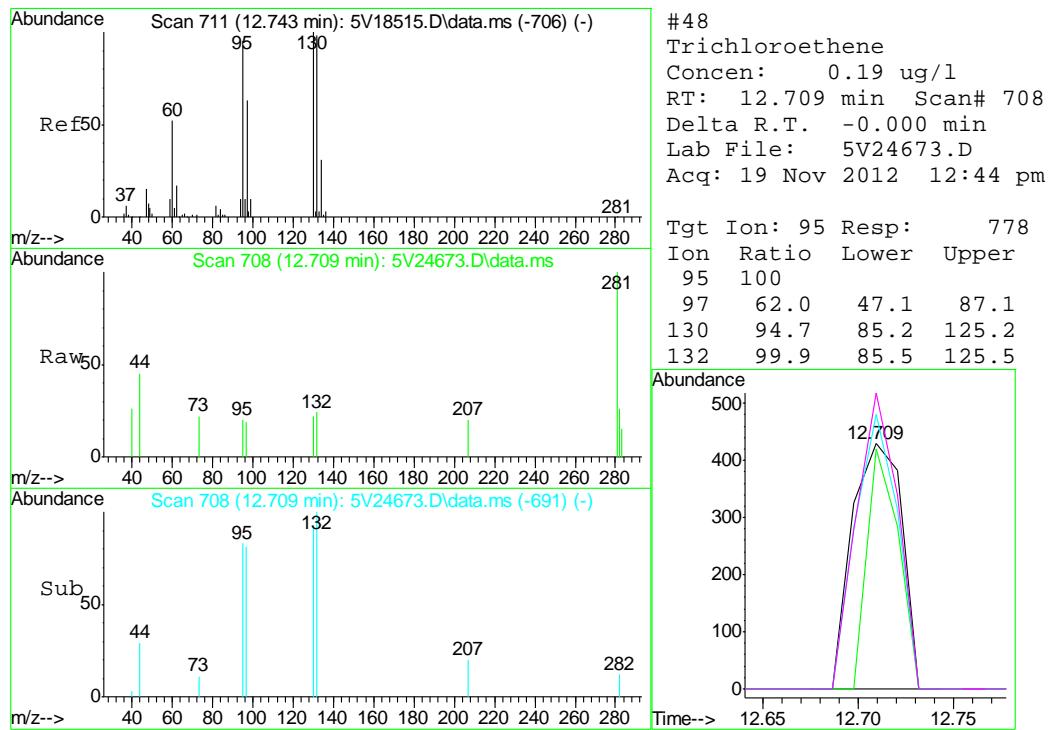
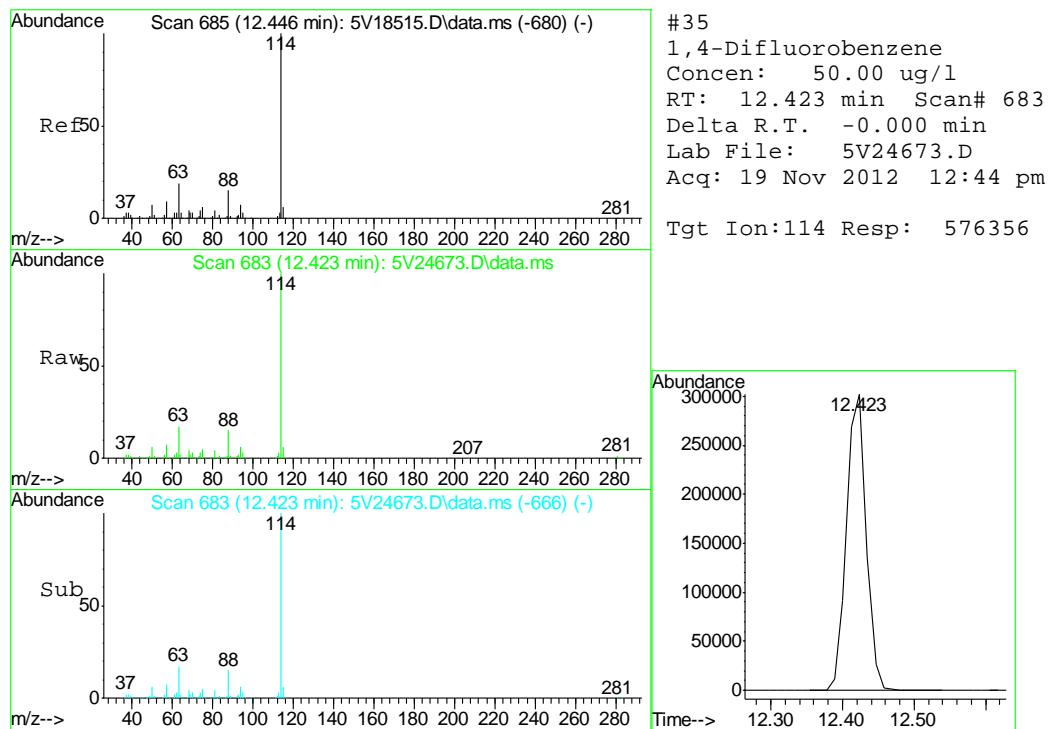
Quantitation Report (QT Reviewed)

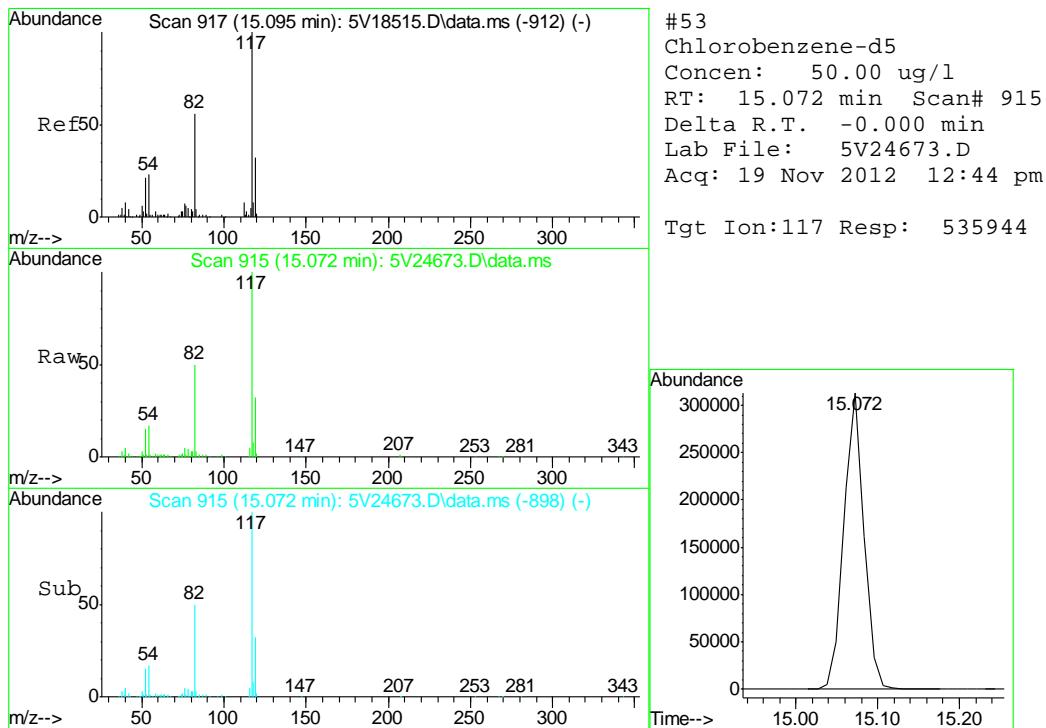
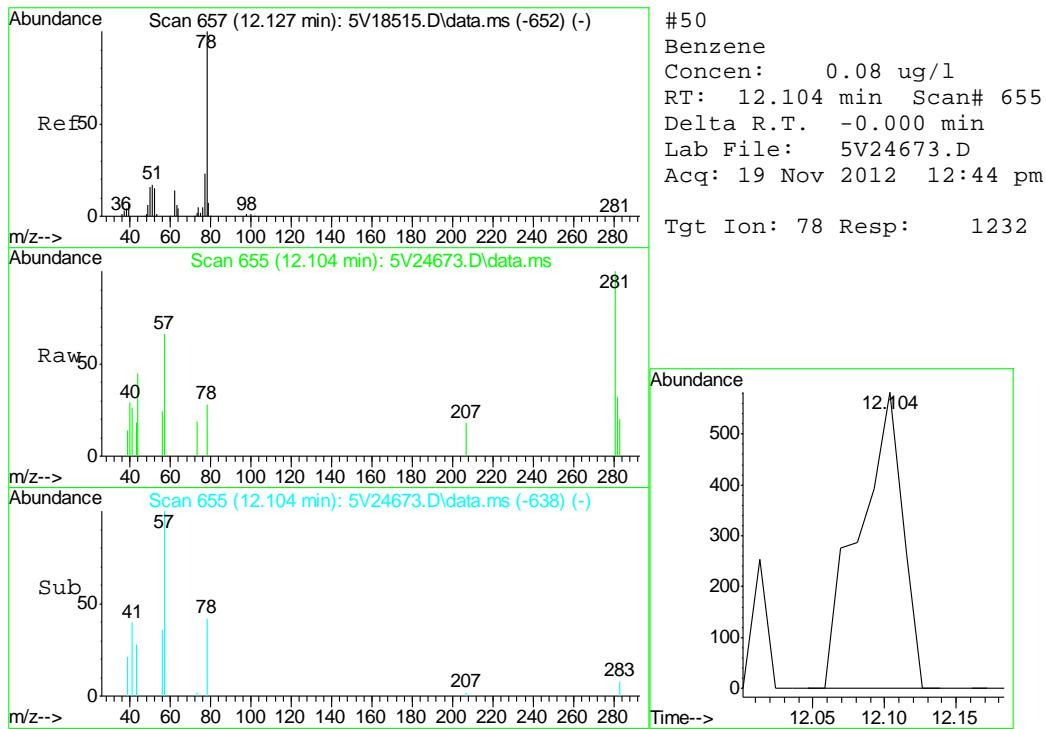
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 Sample : MB
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 ALS Vial : 5 Sample Multiplier: 1

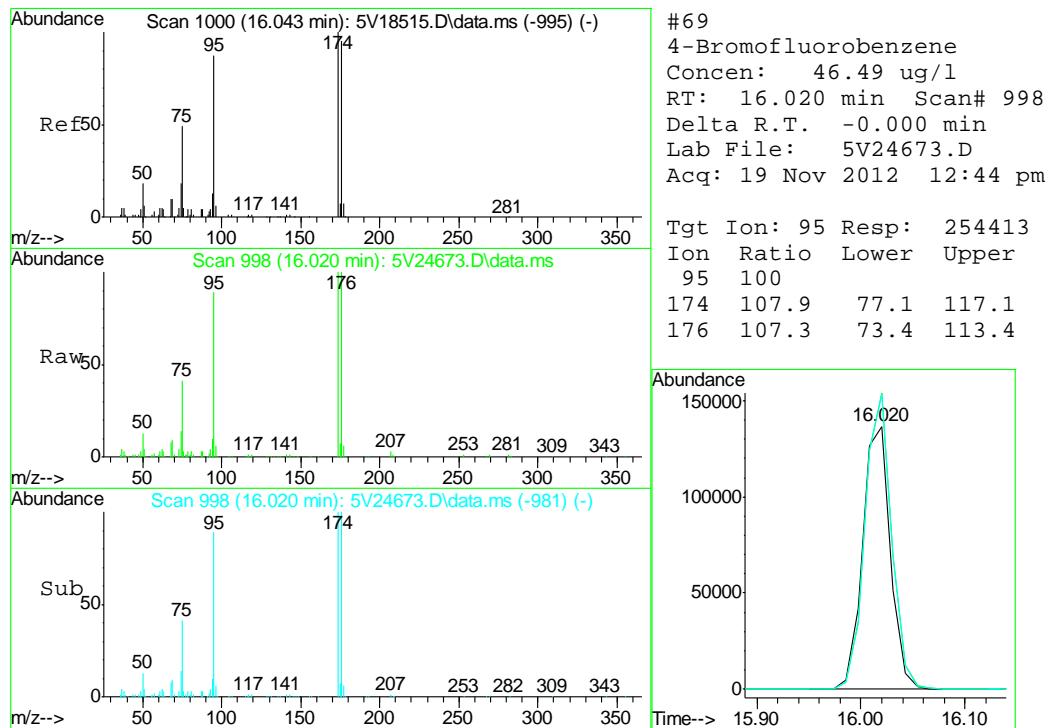
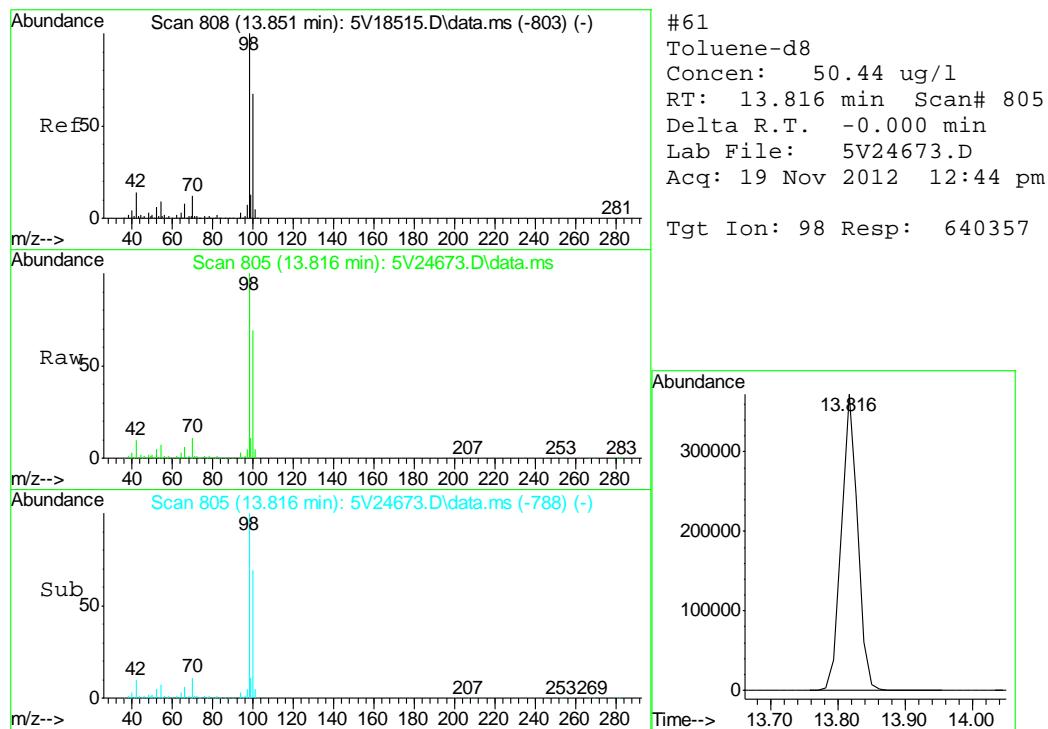
Quant Time: Nov 20 08:43:29 2012
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 Quant Title : 8260
 QLast Update : Wed Nov 14 09:54:38 2012
 Response via : Initial Calibration

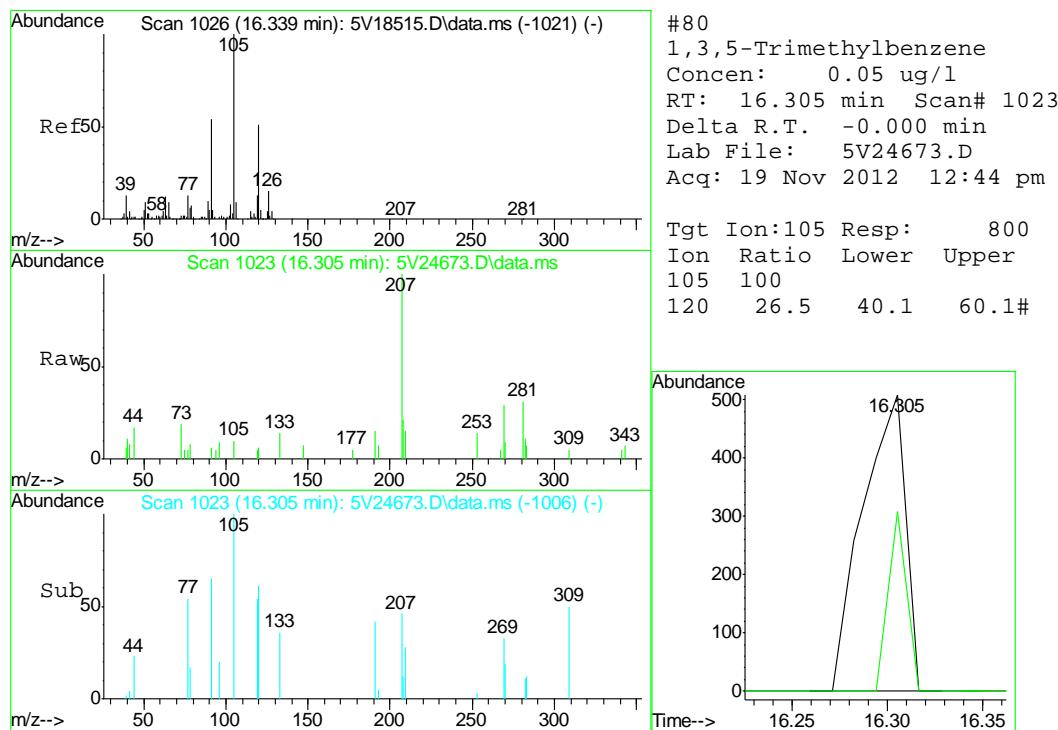
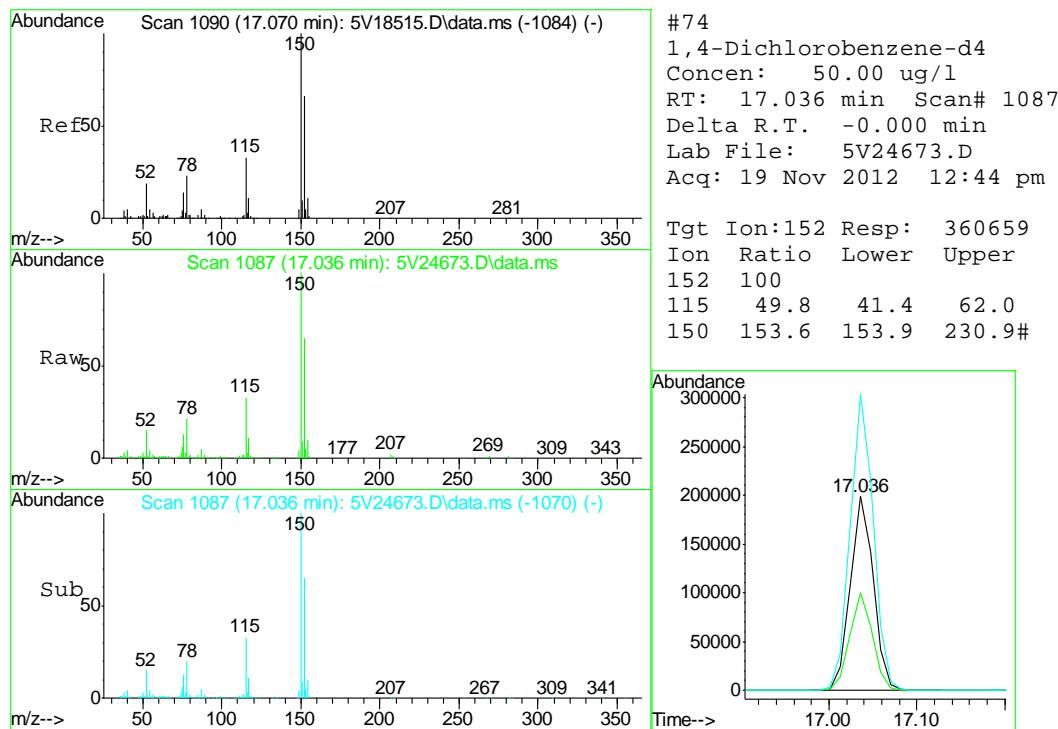


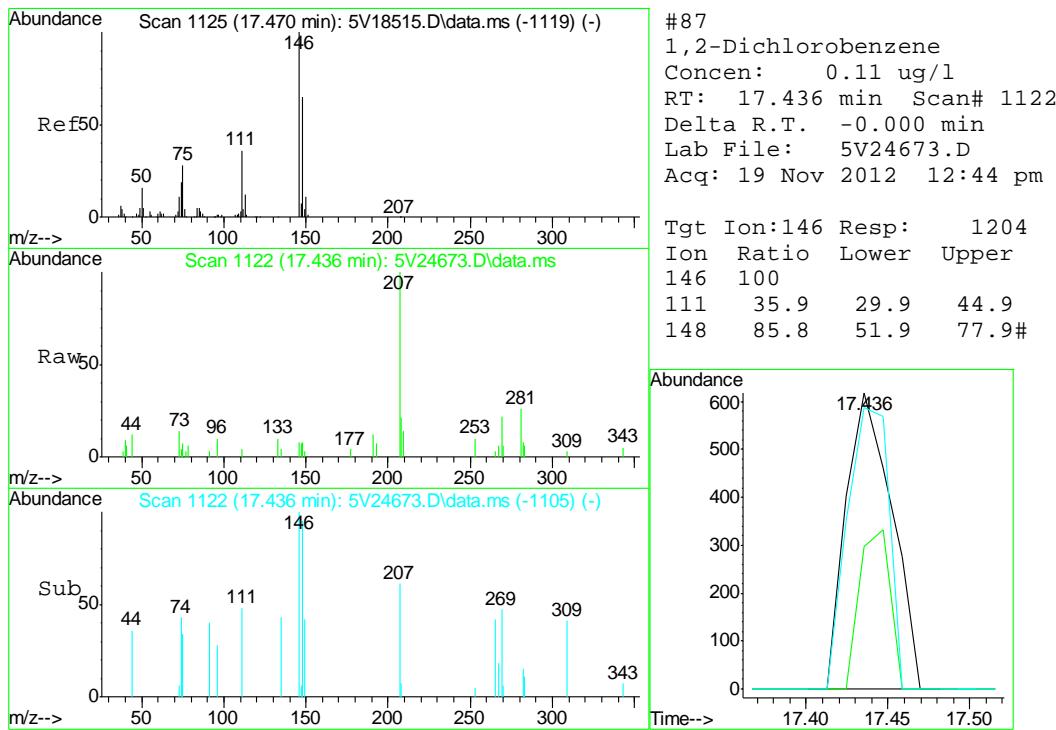
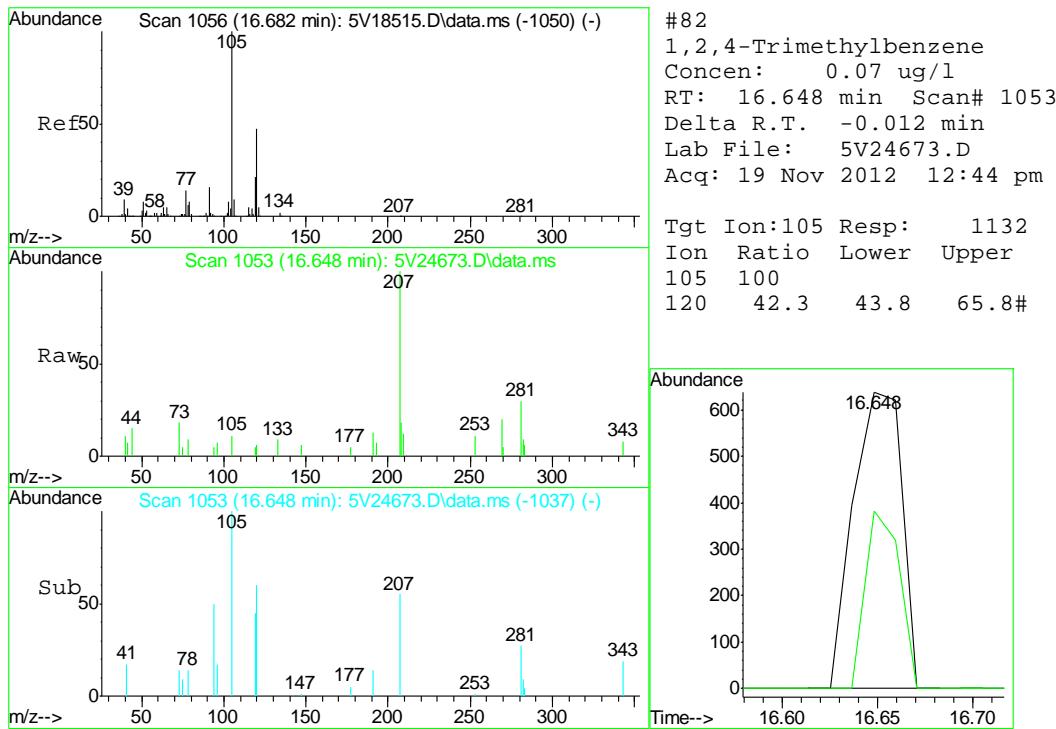


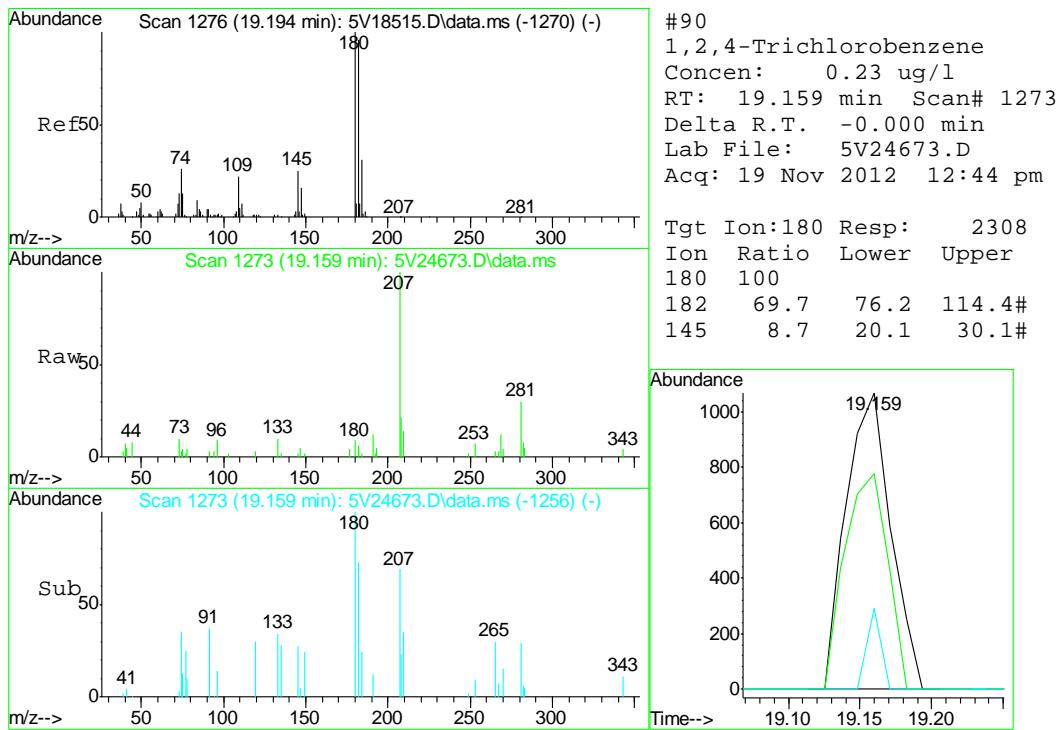
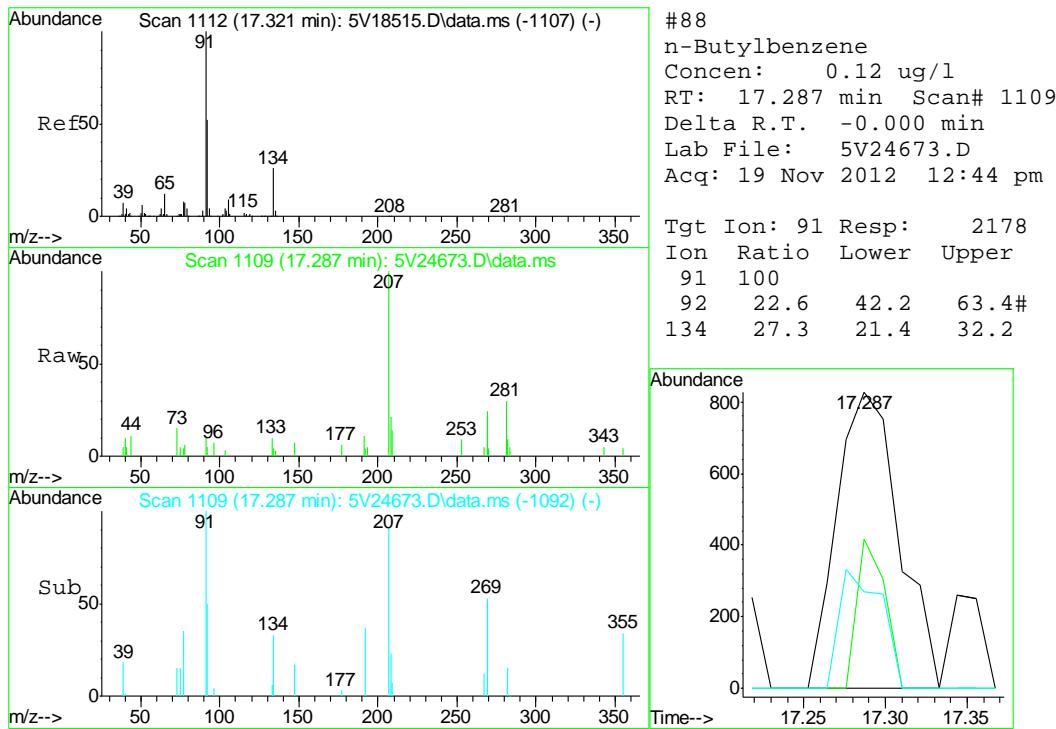


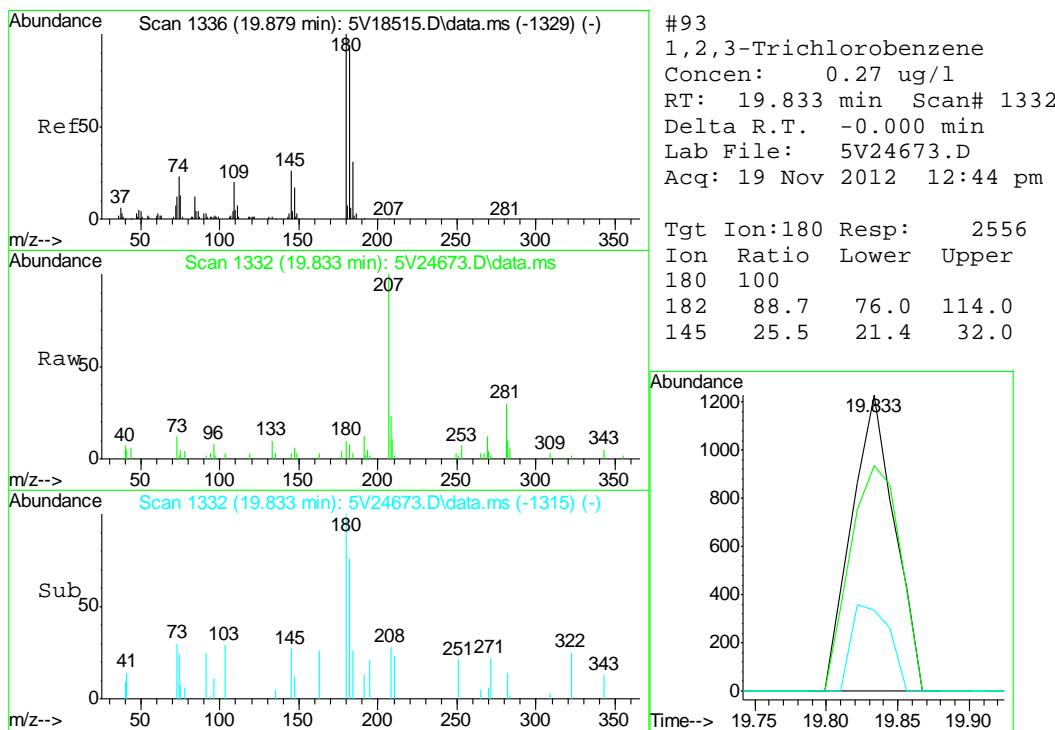
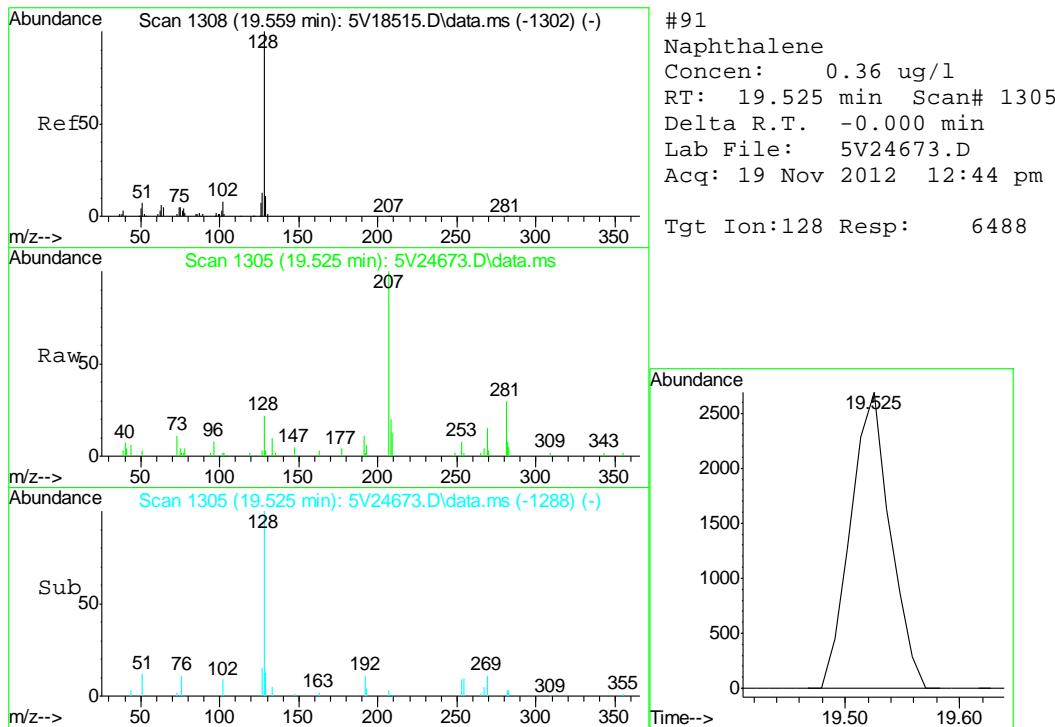














GC/MS Semi-volatiles

QC Data Summaries

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Includes the following where applicable:

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

Method Blank Summary

Job Number: D41014

Account: XTOKWR XTO Energy

Project: PCU 296-5A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
OP6988-MB	3G12229.D	1	11/26/12	SM	11/20/12	OP6988	E3G577

The QC reported here applies to the following samples:

Method: SW846 8270C BY SIM

D41014-1

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	97%	10-159%
321-60-8	2-Fluorobiphenyl	99%	19-131%
1718-51-0	Terphenyl-d14	91%	18-150%

Blank Spike Summary

Page 1 of 1

Job Number: D41014
Account: XTOKWR XTO Energy
Project: PCU 296-5A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
OP6988-BS	3G12230.D	1	11/26/12	SM	11/20/12	OP6988	E3G577

The QC reported here applies to the following samples:

Method: SW846 8270C BY SIM

D41014-1

CAS No.	Compound	Spike ug/kg	BSP ug/kg	BSP %	Limits
83-32-9	Acenaphthene	83.3	93.9	113	68-130
120-12-7	Anthracene	83.3	80.2	96	67-130
56-55-3	Benzo(a)anthracene	83.3	82.1	99	65-130
205-99-2	Benzo(b)fluoranthene	83.3	71.9	86	44-130
207-08-9	Benzo(k)fluoranthene	83.3	74.2	89	56-131
50-32-8	Benzo(a)pyrene	83.3	77.6	93	62-130
218-01-9	Chrysene	83.3	73.8	89	70-130
53-70-3	Dibenzo(a,h)anthracene	83.3	75.6	91	55-130
206-44-0	Fluoranthene	83.3	78.6	94	70-130
86-73-7	Fluorene	83.3	79.4	95	70-130
193-39-5	Indeno(1,2,3-cd)pyrene	83.3	75.5	91	56-130
91-20-3	Naphthalene	83.3	102	122	70-130
129-00-0	Pyrene	83.3	75.6	91	70-130

CAS No.	Surrogate Recoveries	BSP	Limits
4165-60-0	Nitrobenzene-d5	90%	10-159%
321-60-8	2-Fluorobiphenyl	93%	19-131%
1718-51-0	Terphenyl-d14	86%	18-150%

* = Outside of Control Limits.

Matrix Spike/Matrix Spike Duplicate Summary

Page 1 of 1

Job Number: D41014

Account: XTOKWR XTO Energy

Project: PCU 296-5A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
OP6988-MS	3G12232.D	1	11/26/12	SM	11/20/12	OP6988	E3G577
OP6988-MSD	3G12233.D	1	11/26/12	SM	11/20/12	OP6988	E3G577
D41014-1	3G12231.D	1	11/26/12	SM	11/20/12	OP6988	E3G577

The QC reported here applies to the following samples:

Method: SW846 8270C BY SIM

D41014-1

CAS No.	Compound	D41014-1		Spike	MS	MS	MSD	MSD	RPD	Limits Rec/RPD
		ug/kg	Q	ug/kg	ug/kg	%	ug/kg	%		
83-32-9	Acenaphthene	ND		96.2	110	114	103	107	7	25-151/30
120-12-7	Anthracene	ND		96.2	96.8	101	93.9	98	3	39-159/30
56-55-3	Benzo(a)anthracene	ND		96.2	98.2	102	95.3	99	3	39-168/30
205-99-2	Benzo(b)fluoranthene	ND		96.2	83.8	87	83.2	86	1	24-163/30
207-08-9	Benzo(k)fluoranthene	ND		96.2	88.6	92	85.5	89	4	10-188/30
50-32-8	Benzo(a)pyrene	ND		96.2	92.1	96	88.2	92	4	32-144/30
218-01-9	Chrysene	ND		96.2	88.5	92	84.3	88	5	43-150/30
53-70-3	Dibenzo(a,h)anthracene	ND		96.2	84.5	88	85.3	89	1	21-152/30
206-44-0	Fluoranthene	ND		96.2	99.2	103	95.6	99	4	36-157/30
86-73-7	Fluorene	ND		96.2	98.0	102	94.0	98	4	10-182/30
193-39-5	Indeno(1,2,3-cd)pyrene	ND		96.2	87.0	90	84.6	88	3	20-154/30
91-20-3	Naphthalene	34.8		96.2	134	103	111	79	19	10-163/30
129-00-0	Pyrene	ND		96.2	95.0	99	91.8	95	3	25-180/30

CAS No.	Surrogate Recoveries	MS	MSD	D41014-1	Limits
4165-60-0	Nitrobenzene-d5	86%	74%	80%	10-159%
321-60-8	2-Fluorobiphenyl	84%	80%	85%	19-131%
1718-51-0	Terphenyl-d14	80%	78%	85%	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\112612\
 Data File : 3g12231.D
 Acq On : 26 Nov 2012 6:54 pm
 Operator : SARAHM1
 Sample : D41014-1
 Misc : OP6988,E3G577,30.04,,,1,1
 ALS Vial : 9 Sample Multiplier: 1

Quant Time: Nov 27 09:24:41 2012
 Quant Method : C:\msdchem\1\METHODS\SIMPE3G574.M
 Quant Title : PAHSIM BASE
 QLast Update : Mon Nov 26 15:39:31 2012
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Naphthalene-d8	5.707	136	137587	4.0000	ug/mL	0.00
6) Acenaphthene-d10	7.412	164	88246	4.0000	ug/mL	0.00
15) Phenanthrene-d10	8.894	188	156268	4.0000	ug/mL	-0.06
19) Chrysene-d12	11.535	240	124910	4.0000	ug/mL	0.00
24) Perylene-d12	12.913	264	73114	4.0000	ug/mL	0.00

System Monitoring Compounds

2) Nitrobenzene-d5	5.021	82	531547	40.1835	ug/mL	0.00
Spiked Amount	50.000	Range	25 - 135	Recovery	=	80.36%
7) 2-Fluorobiphenyl	6.751	172	1363494	42.3736	ug/mL	0.00
Spiked Amount	50.000	Range	25 - 135	Recovery	=	84.74%
21) Terphenyl-d14	10.493	244	689124	42.3295	ug/mL	0.00
Spiked Amount	50.000	Range	25 - 135	Recovery	=	84.66%

Target Compounds

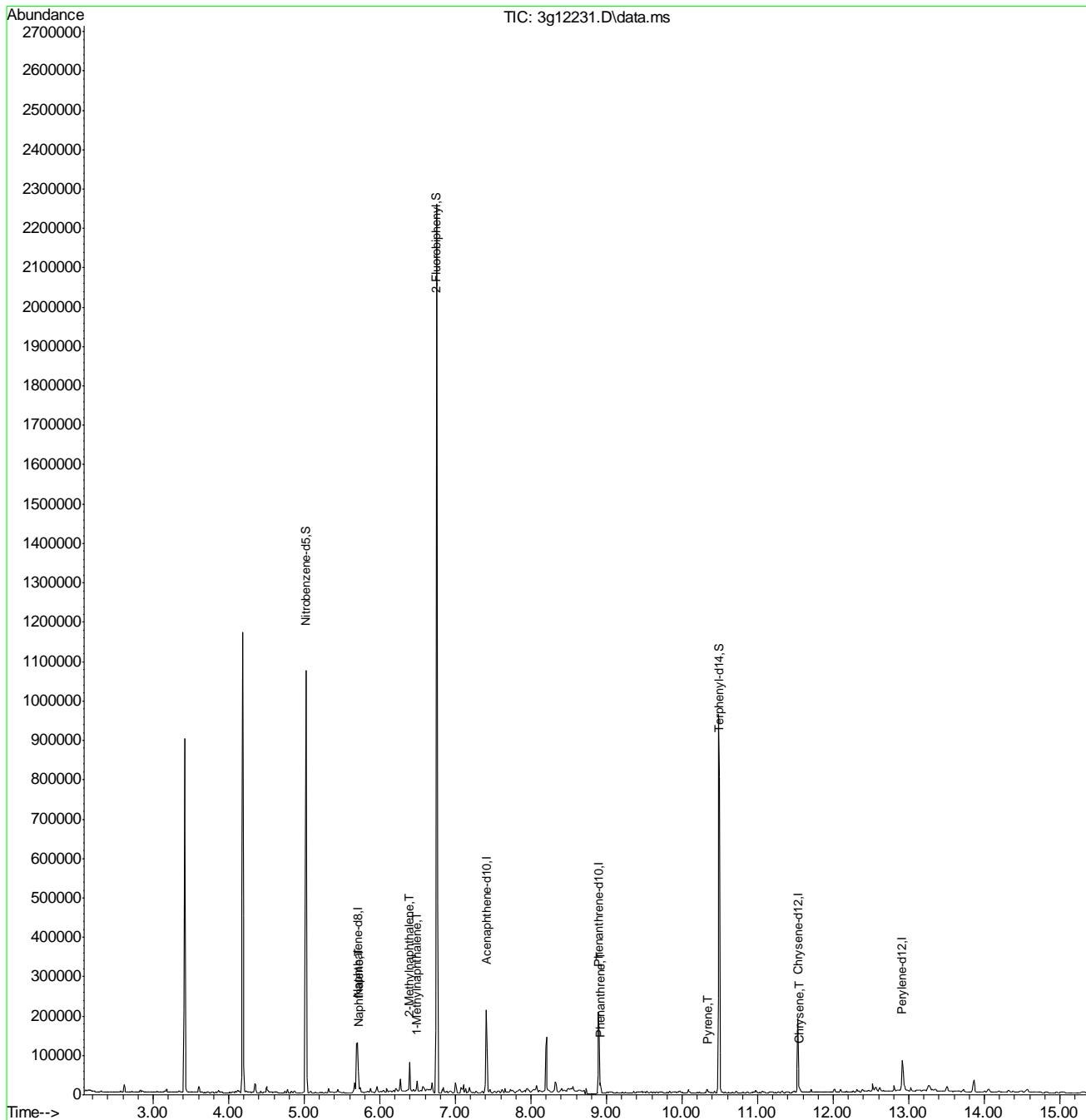
				Qvalue
3) N-Nitrosodimethylamine	2.443	74	129	N.D.
4) N-Nitrosodi-propylamine	0.000	70	0	N.D. d
5) Naphthalene	5.719	128	32496	0.9030 ug/mL 95
8) 2-Methylnaphthalene	6.392	142	29666	1.0768 ug/mL 97
9) 1-Methylnaphthalene	6.492	142	10901	0.4244 ug/mL 90
10) Acenaphthylene	7.271	152	489	N.D.
11) Acenaphthene	7.448	154	415	Below Cal # 1
12) Dibenzofuran	7.625	168	1779	N.D.
13) Fluorene	0.000	166	0	N.D. d
14) Diphenylamine	0.000	169	0	N.D. d
16) Phenanthrene	8.918	178	16041	0.2573 ug/mL 80
17) Anthracene	0.000	178	0	N.D. d
18) Fluoranthene	10.105	202	2970	N.D.
20) Pyrene	10.335	202	4918	0.0717 ug/mL# 66
22) Benzo(a)anthracene	0.000	228	0	N.D. d
23) Chrysene	11.555	228	4302	0.0708 ug/mL 92
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	12.860	252	1575	N.D.
28) Indeno(1,2,3-cd)pyrene	14.143	276	823	N.D.
29) Dibenz(a,h)anthracene	14.164	278	803	N.D.
30) Benzo(g,h,i)perylene	14.880	276	181	N.D.

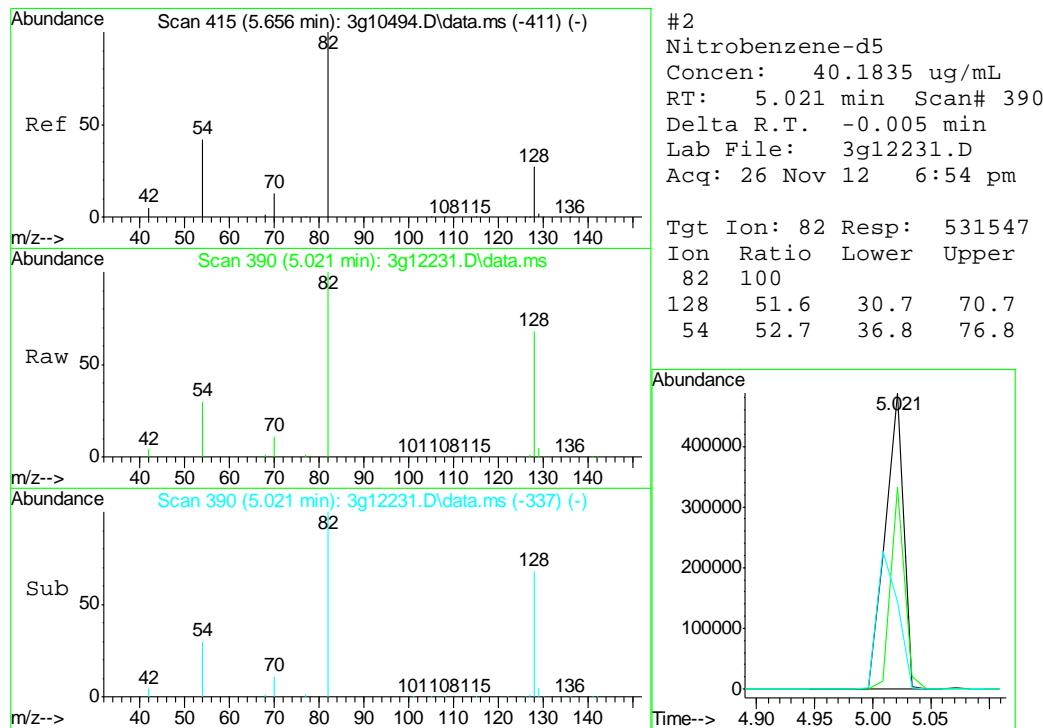
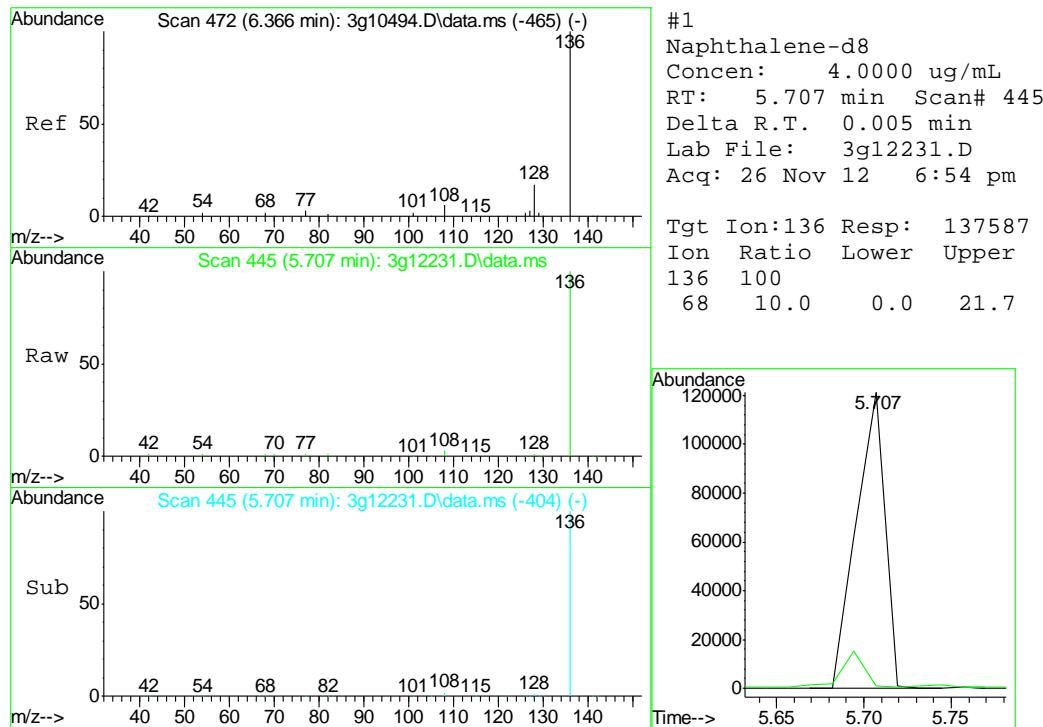
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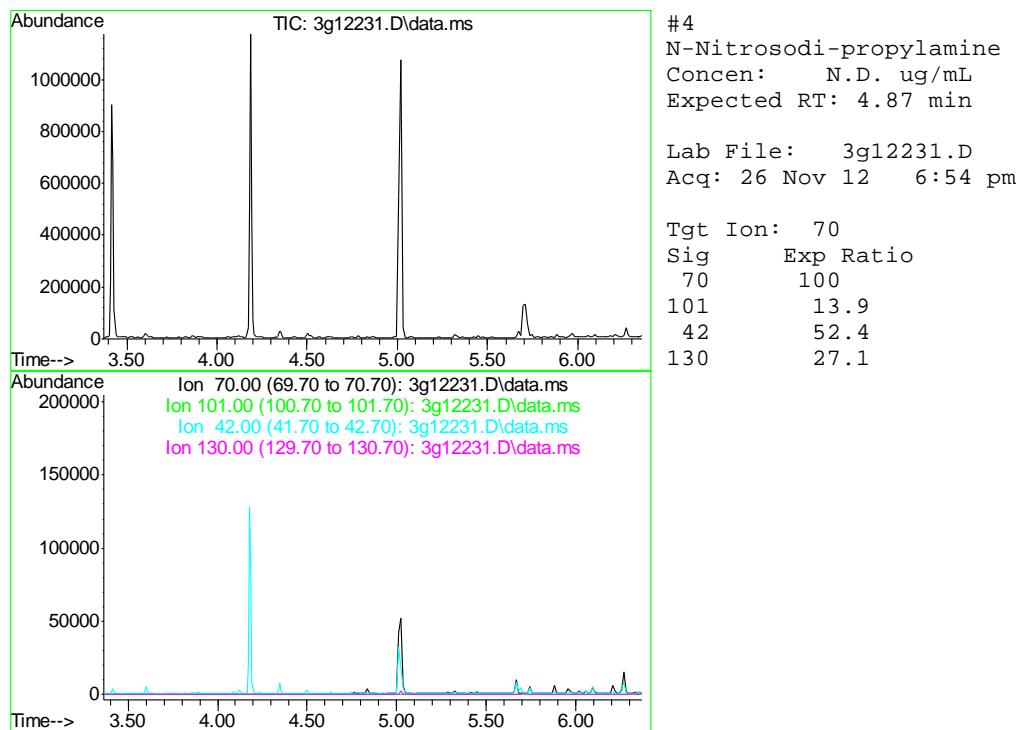
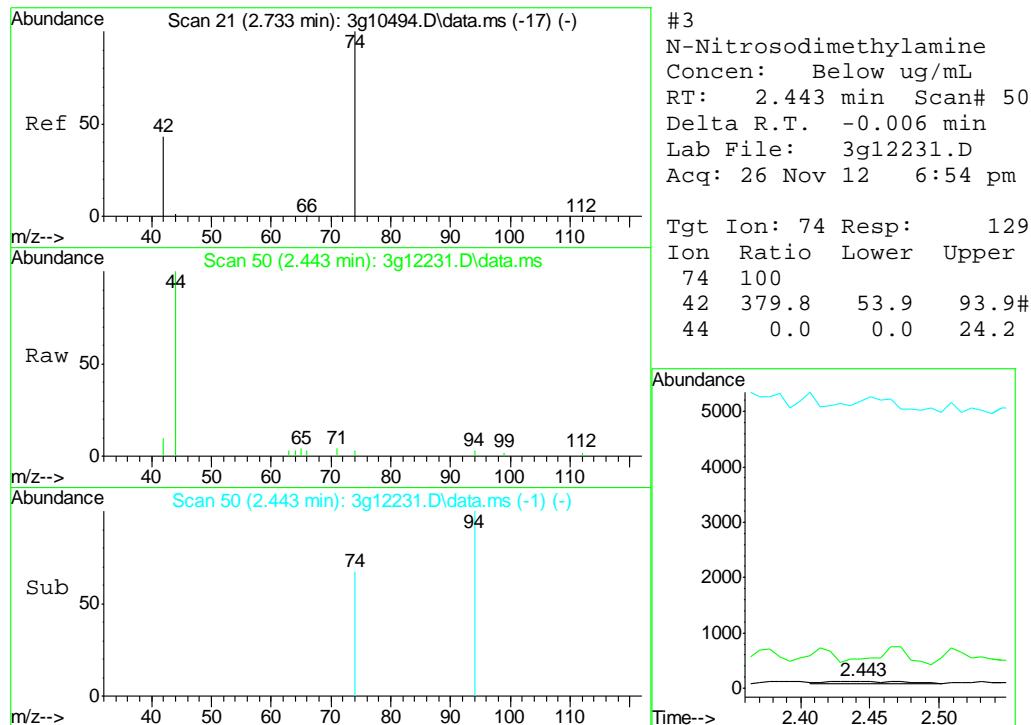
Quantitation Report (QT Reviewed)

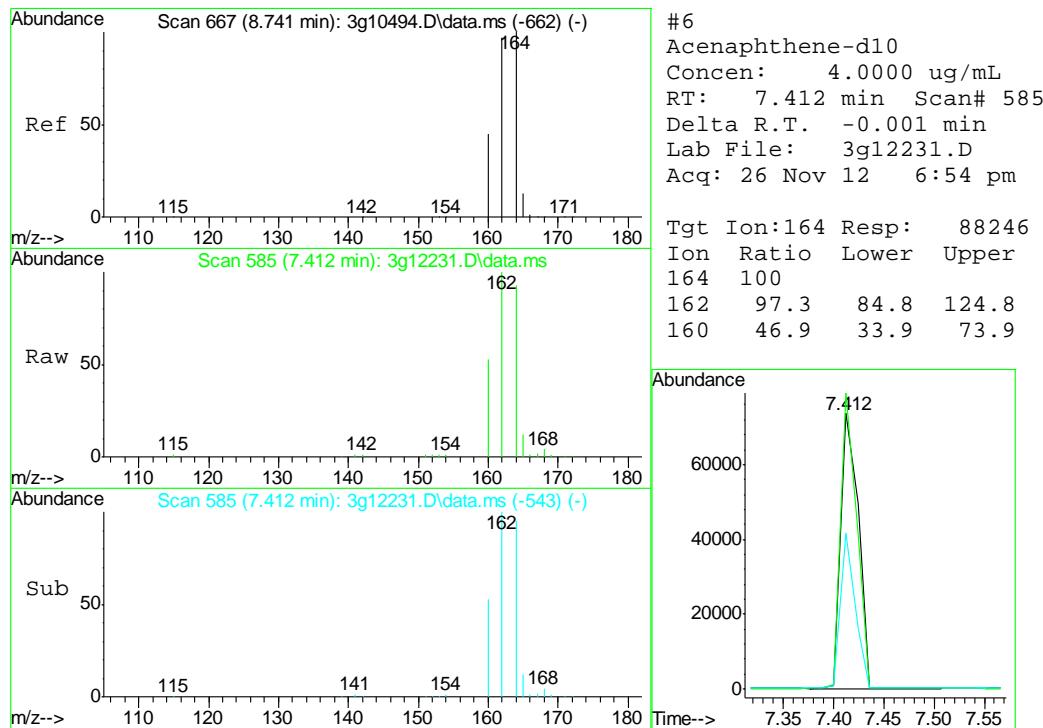
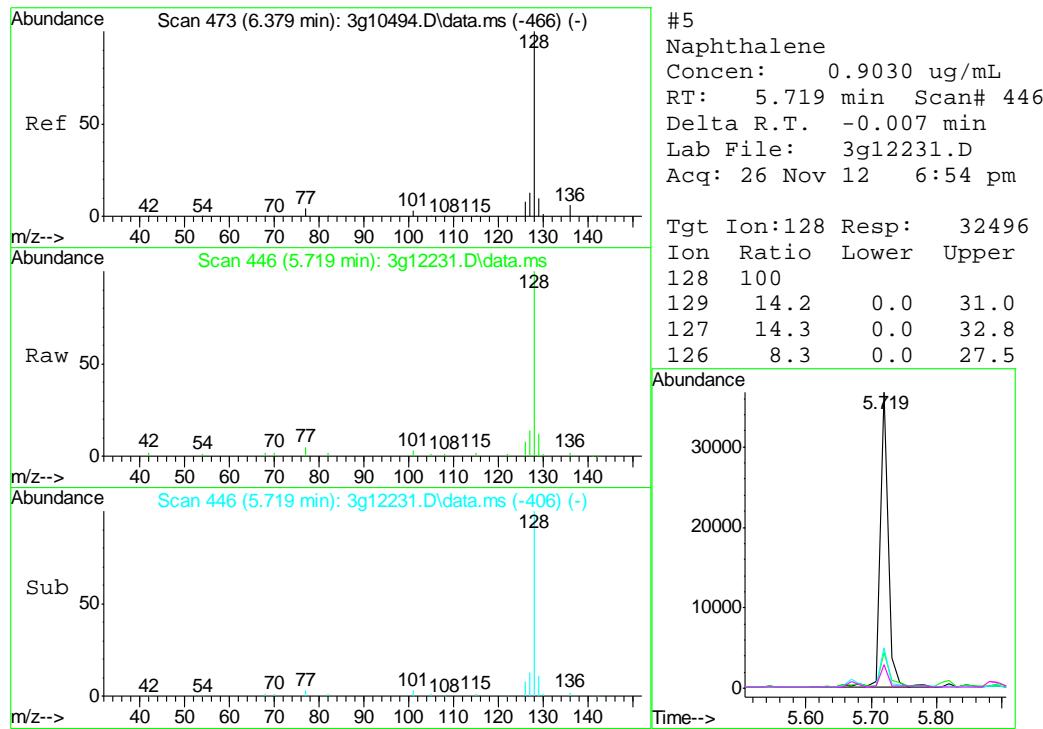
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 Operator : SARAHM1
 Sample : D41014-1
 Misc : OP6988,E3G577,30.04,,,1,1
 ALS Vial : 9 Sample Multiplier: 1

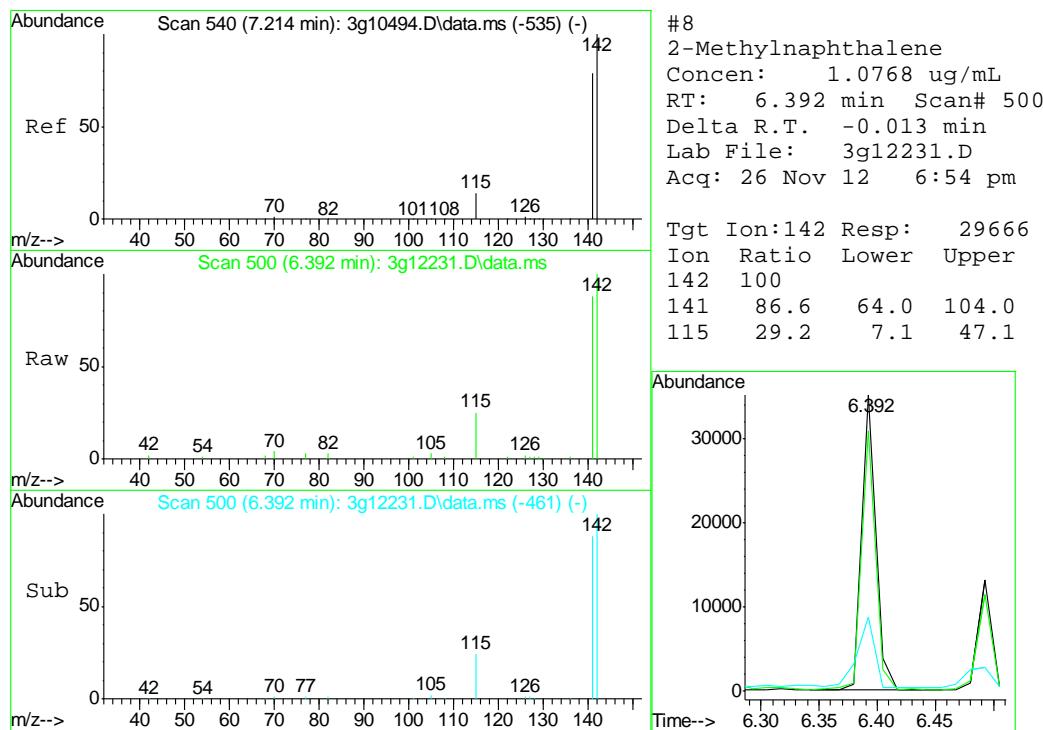
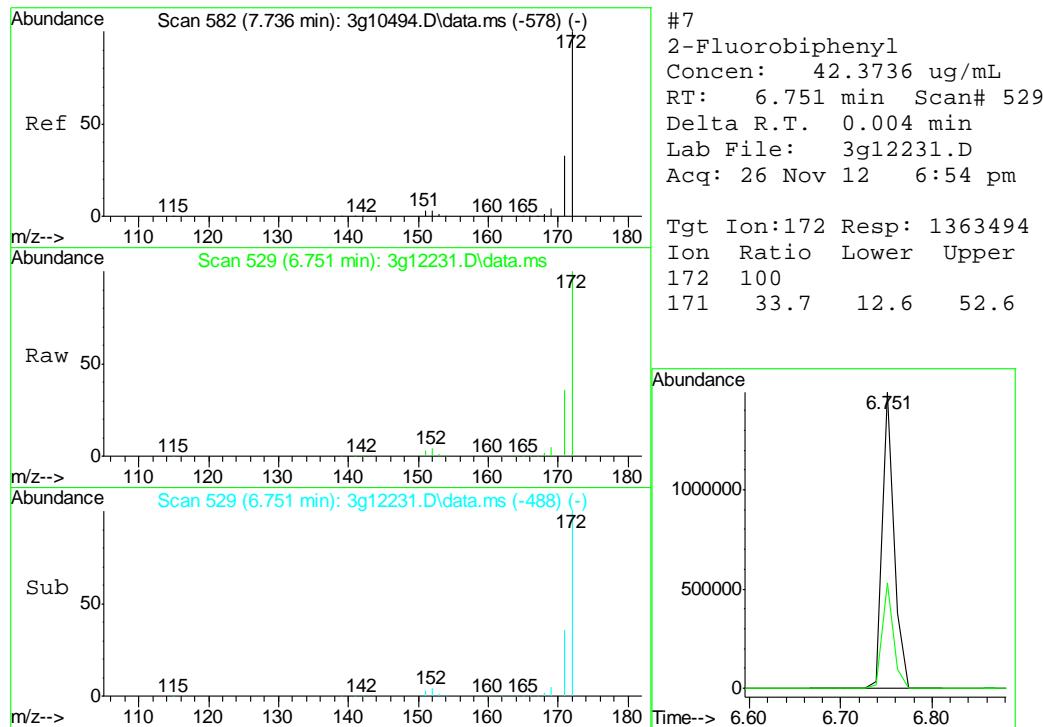
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 Quant Title : PAHSIM BASE
 QLast Update : Mon Nov 26 15:39:31 2012
 Response via : Initial Calibration

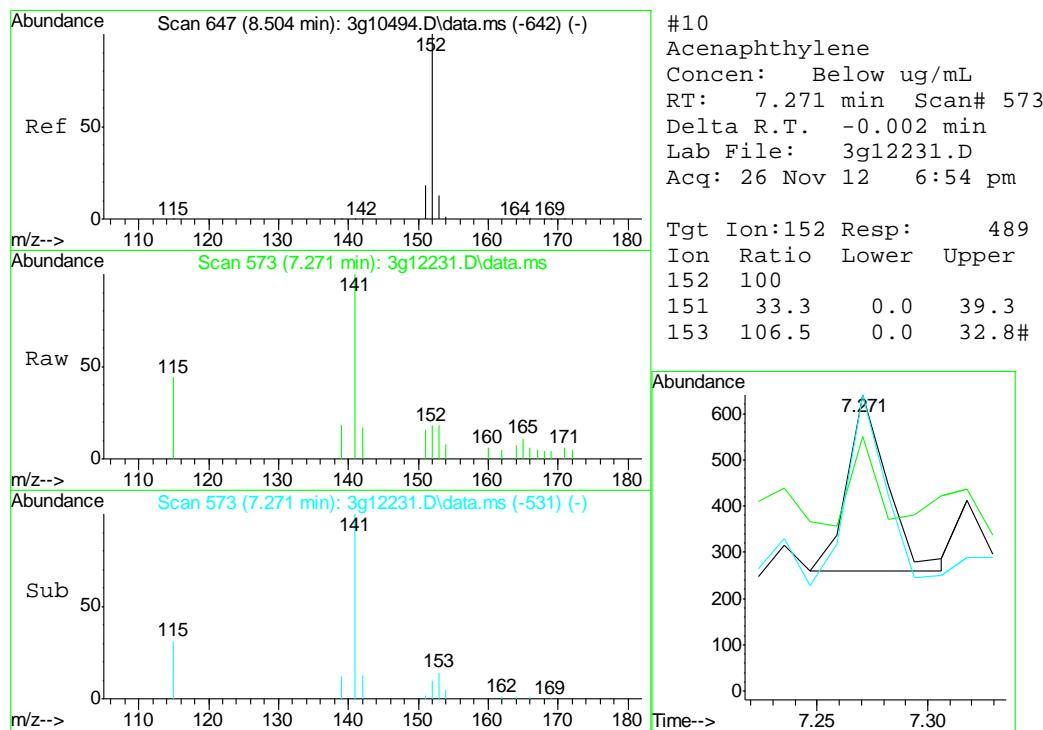
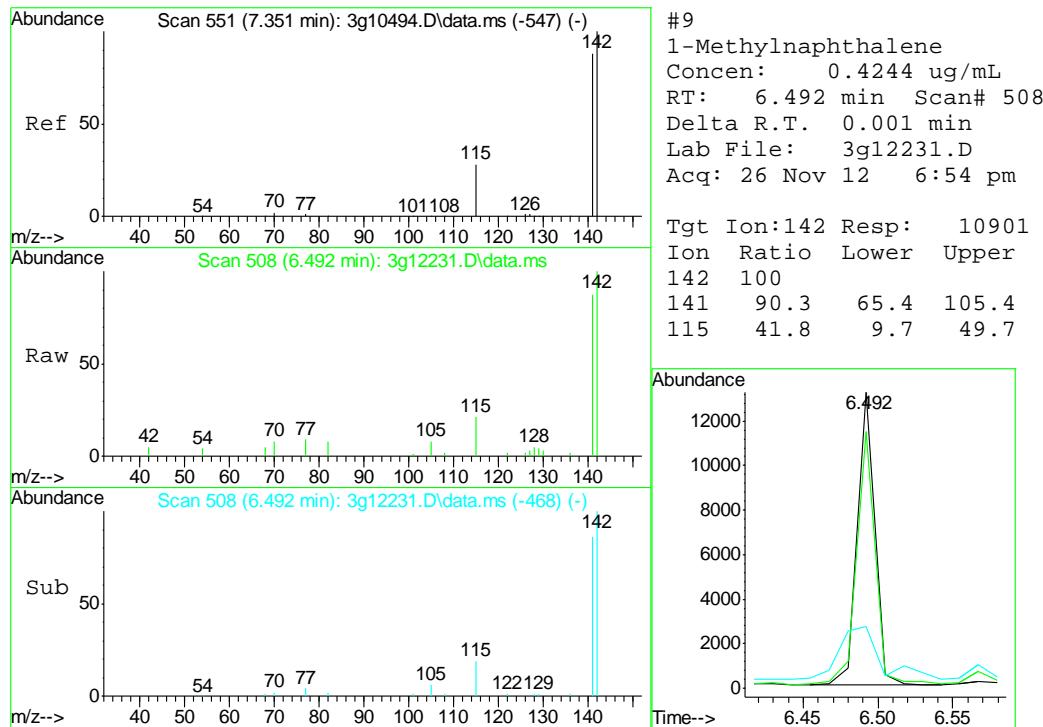


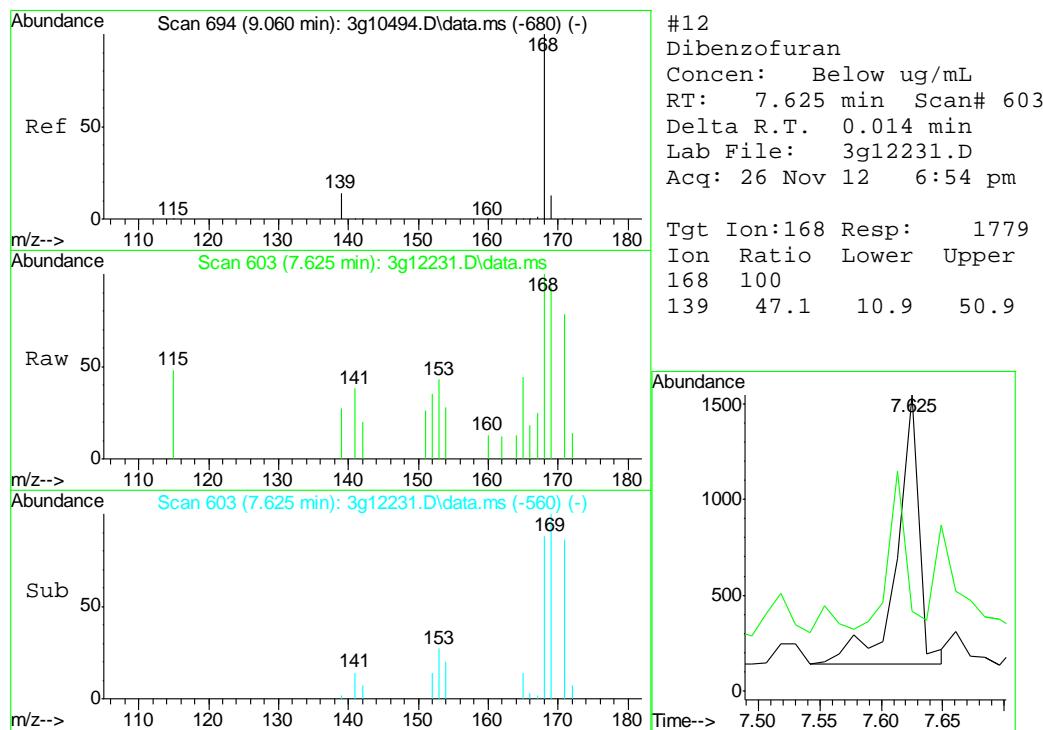
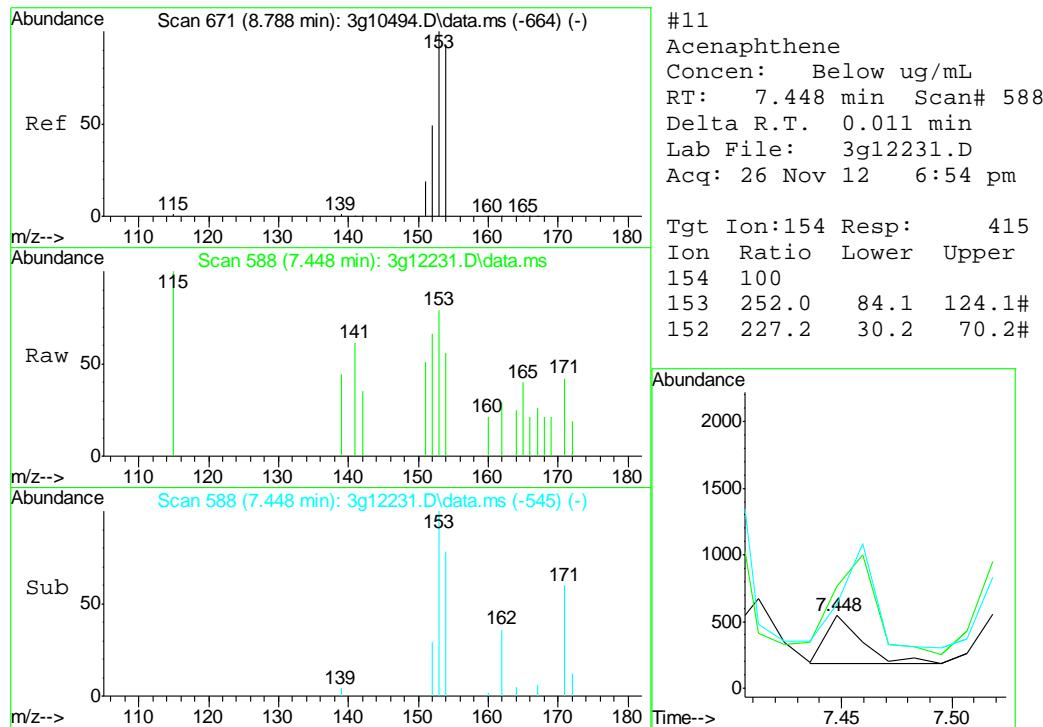


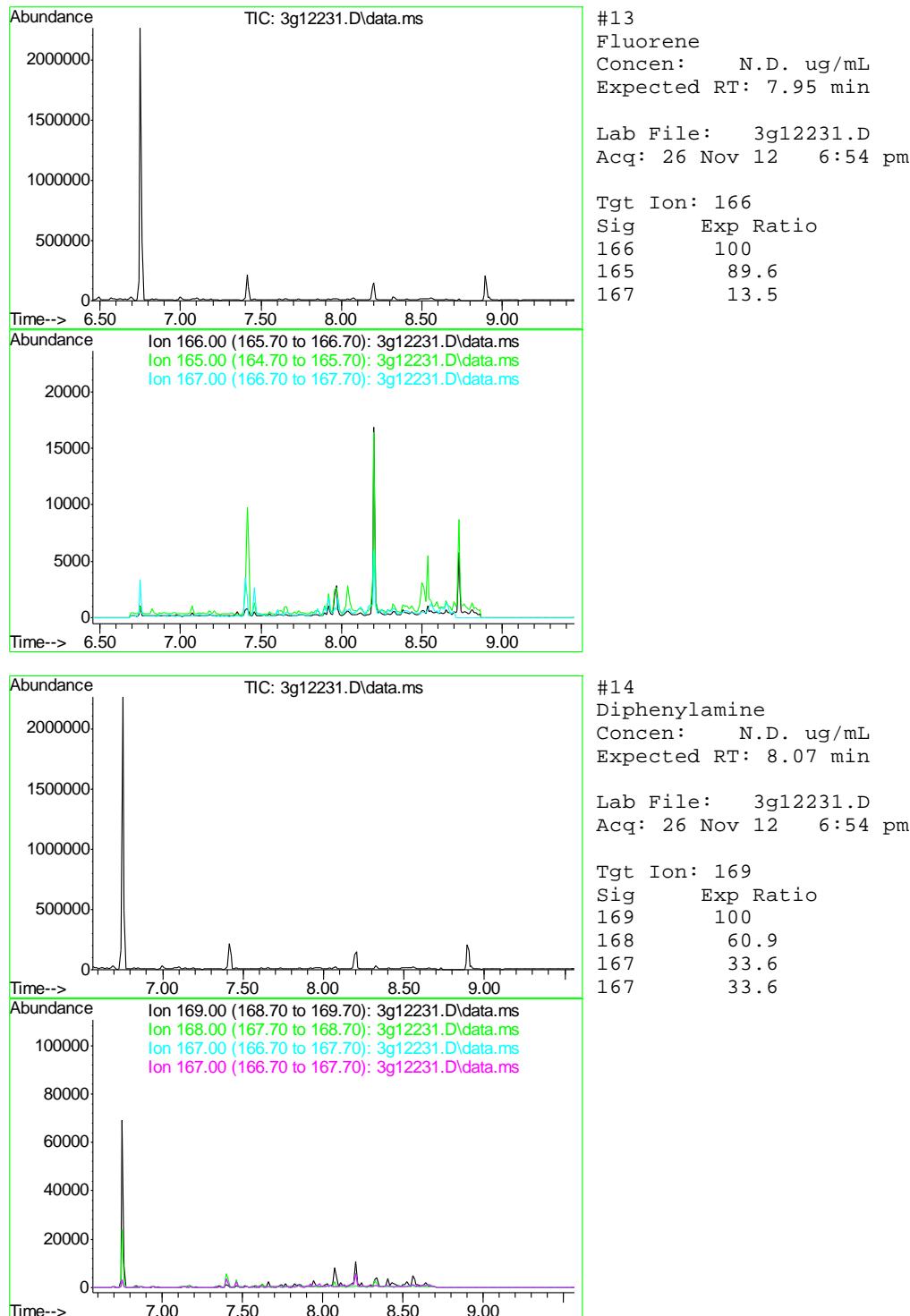


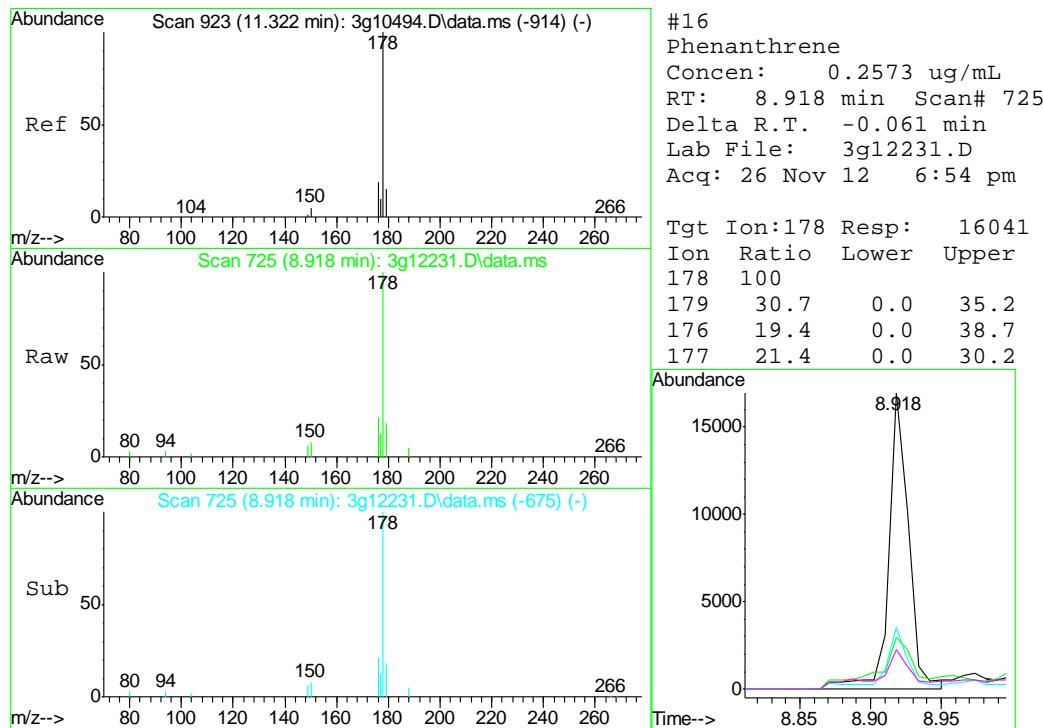
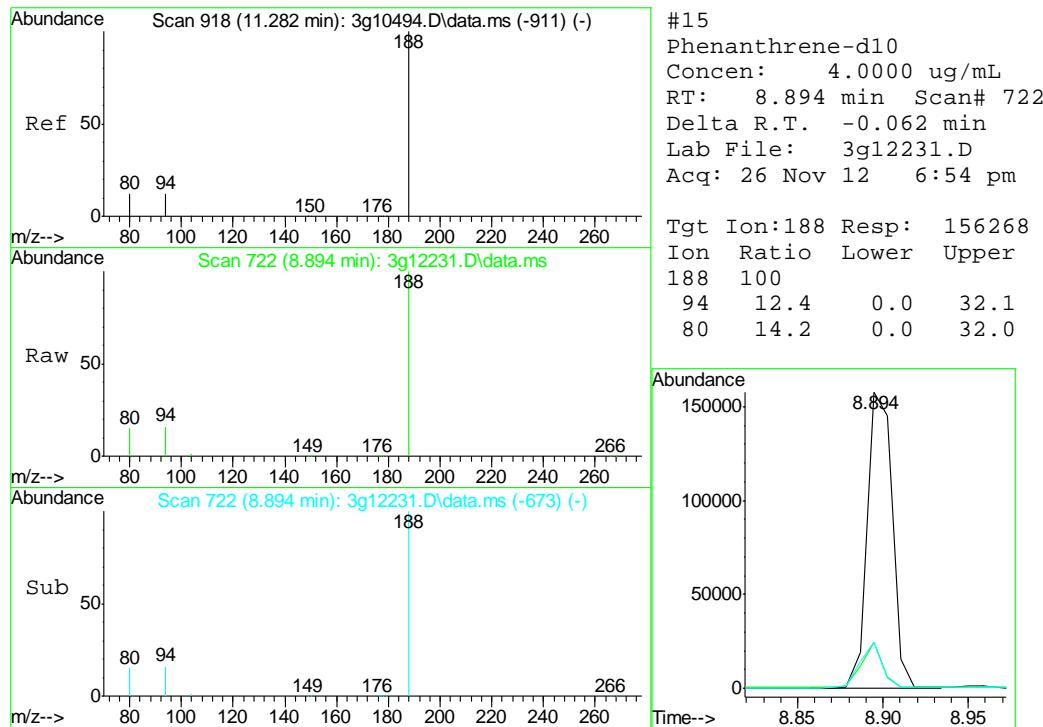


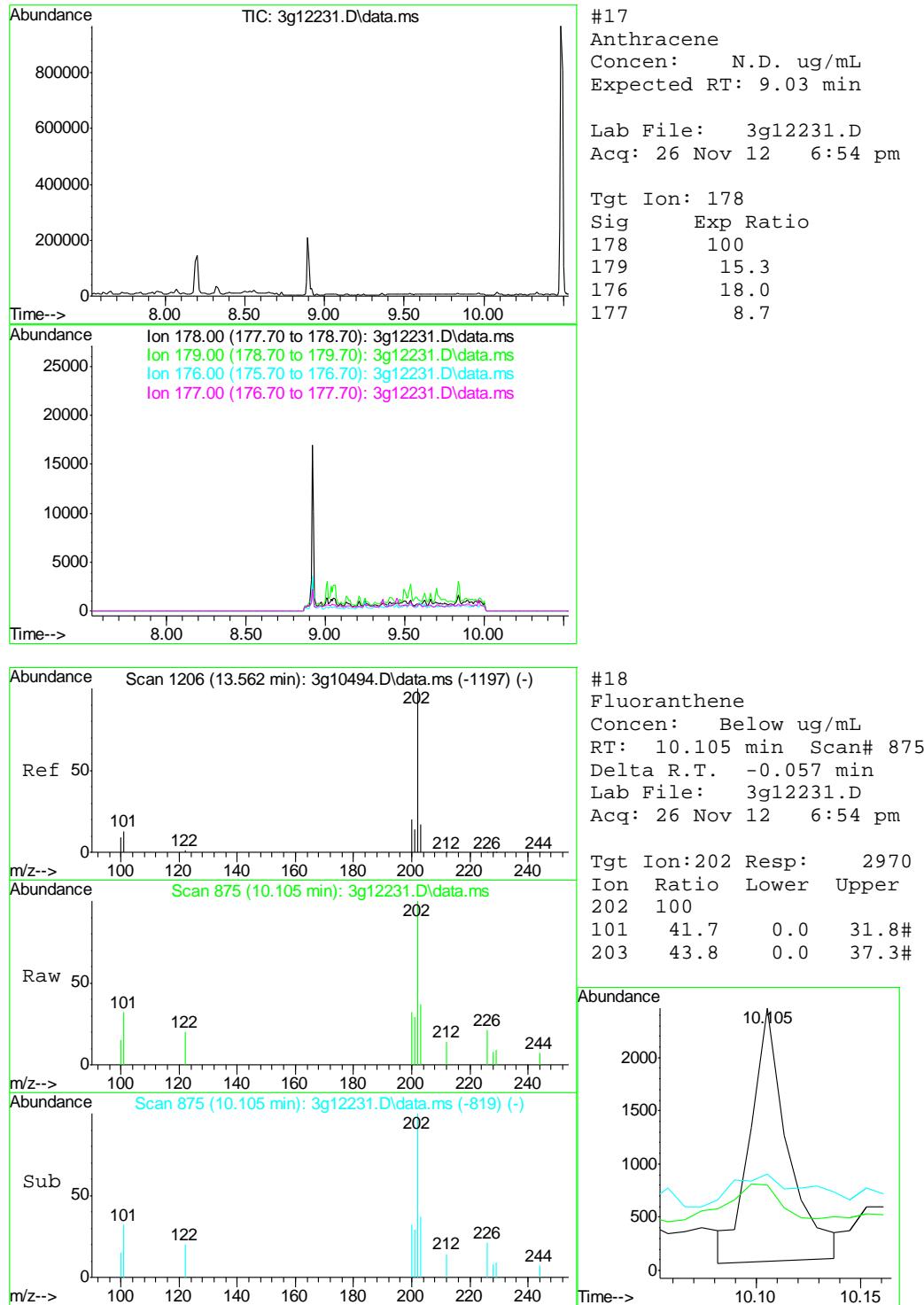


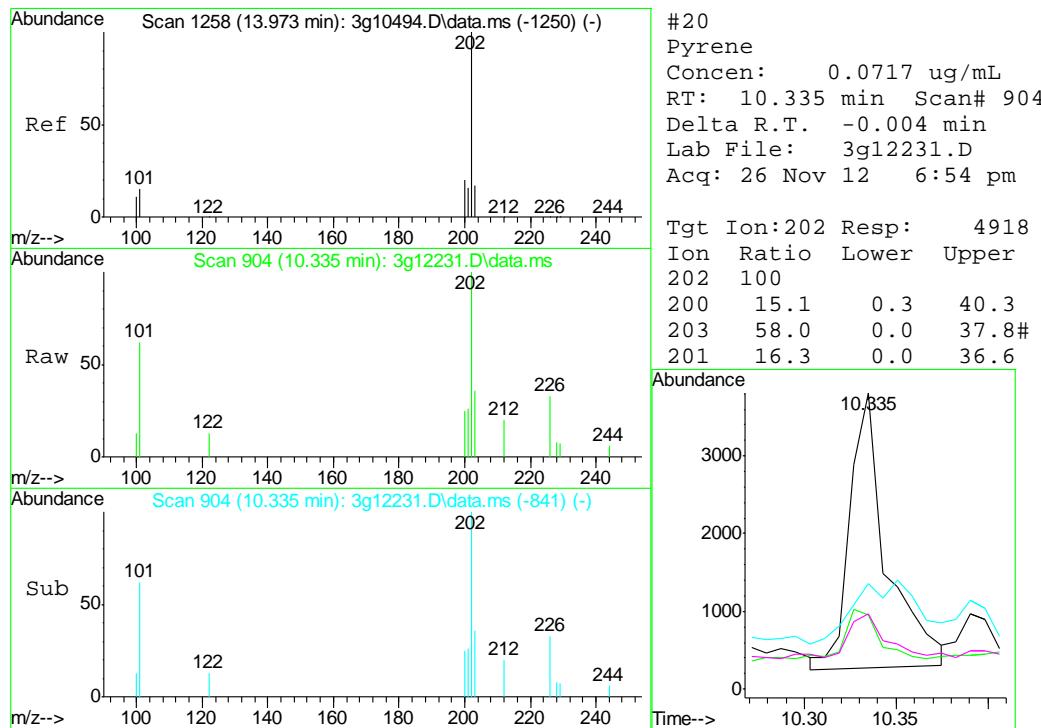
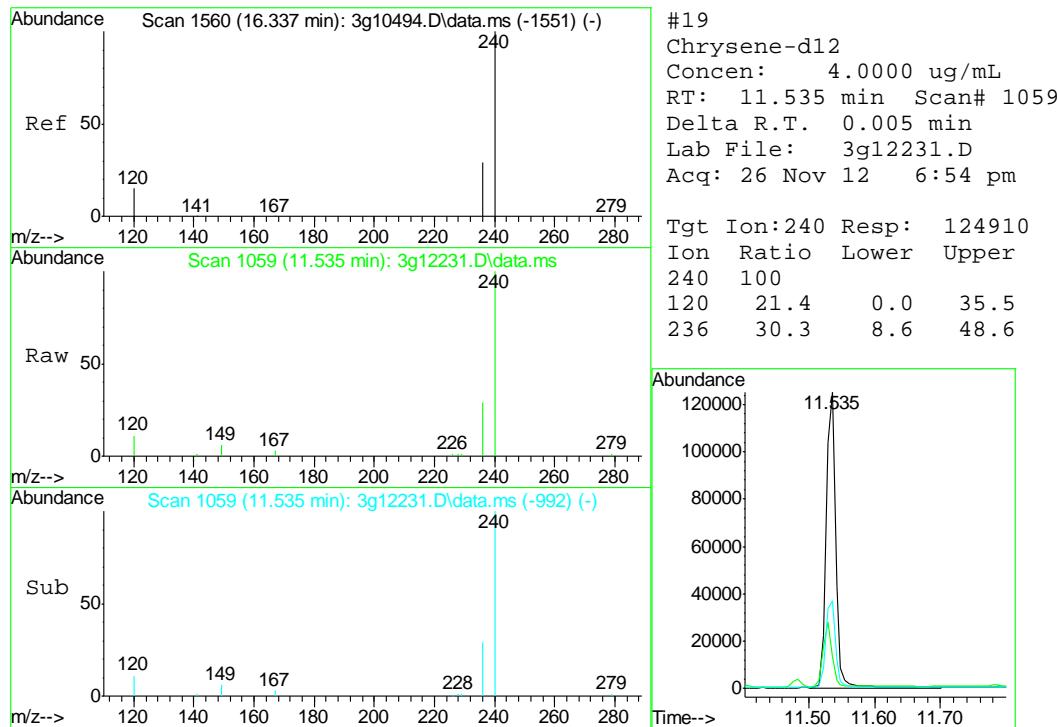


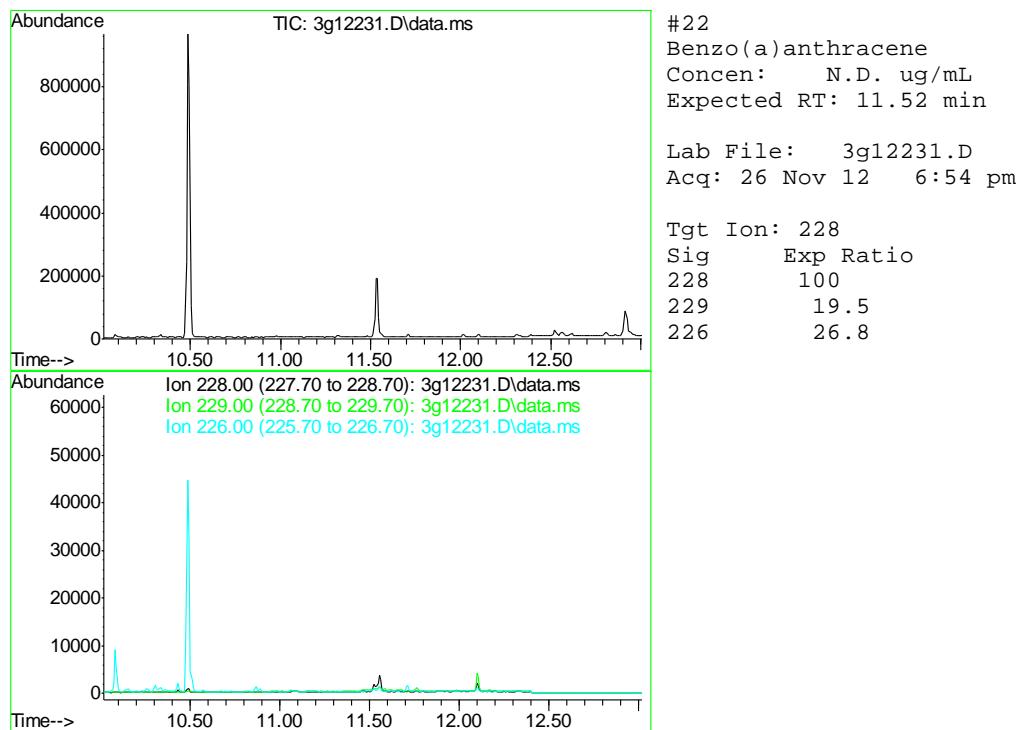
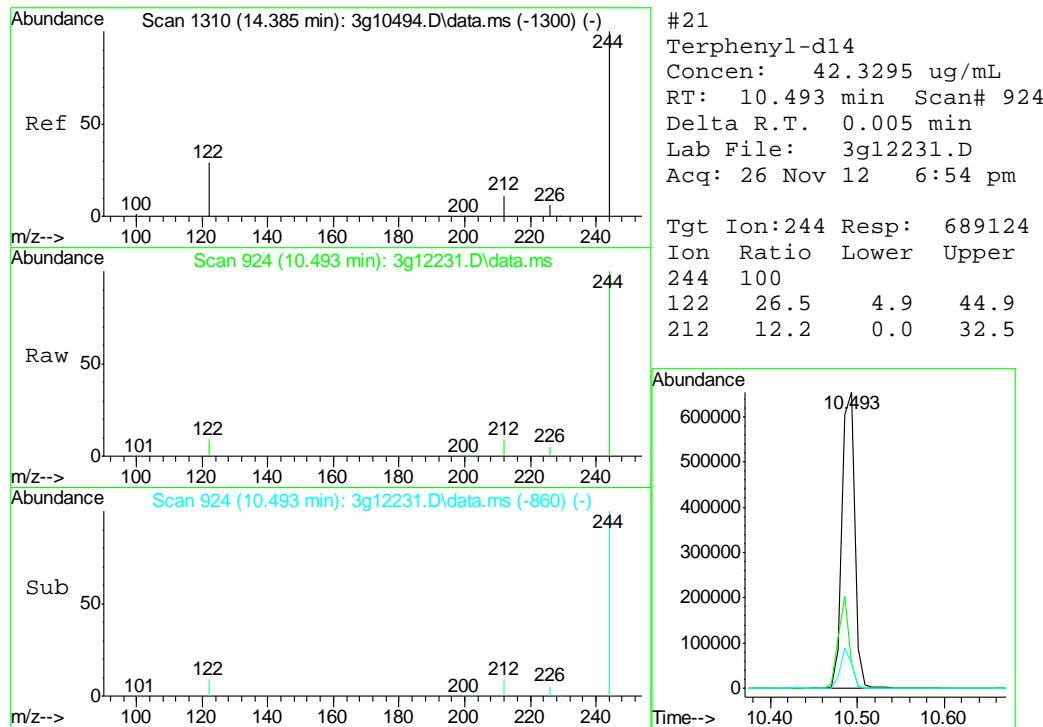


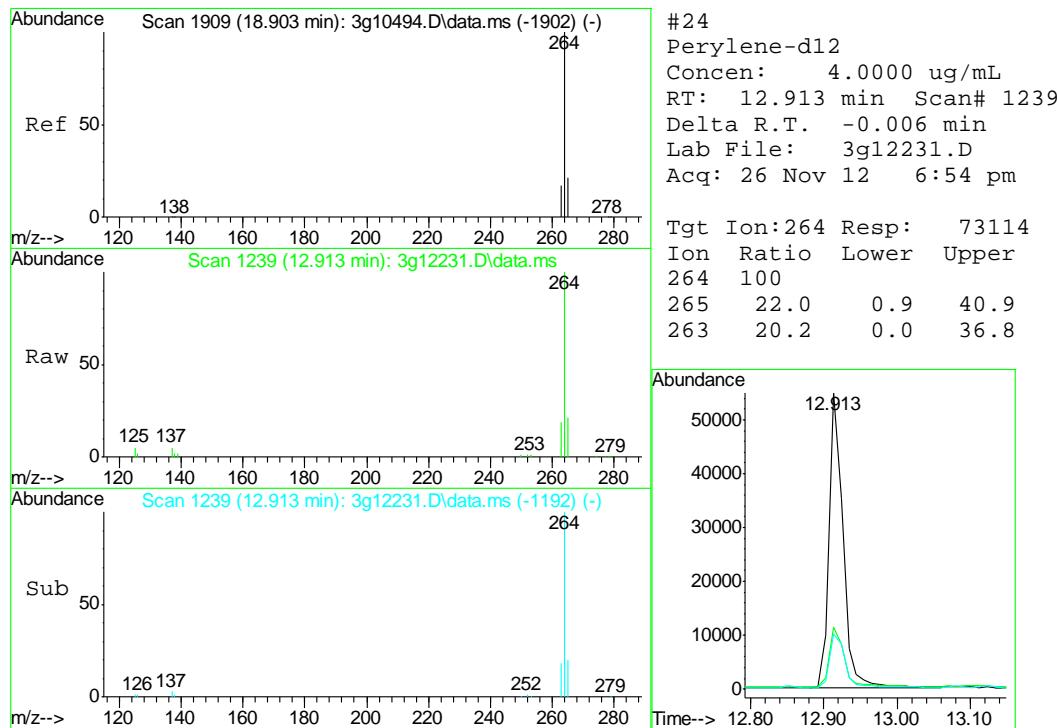
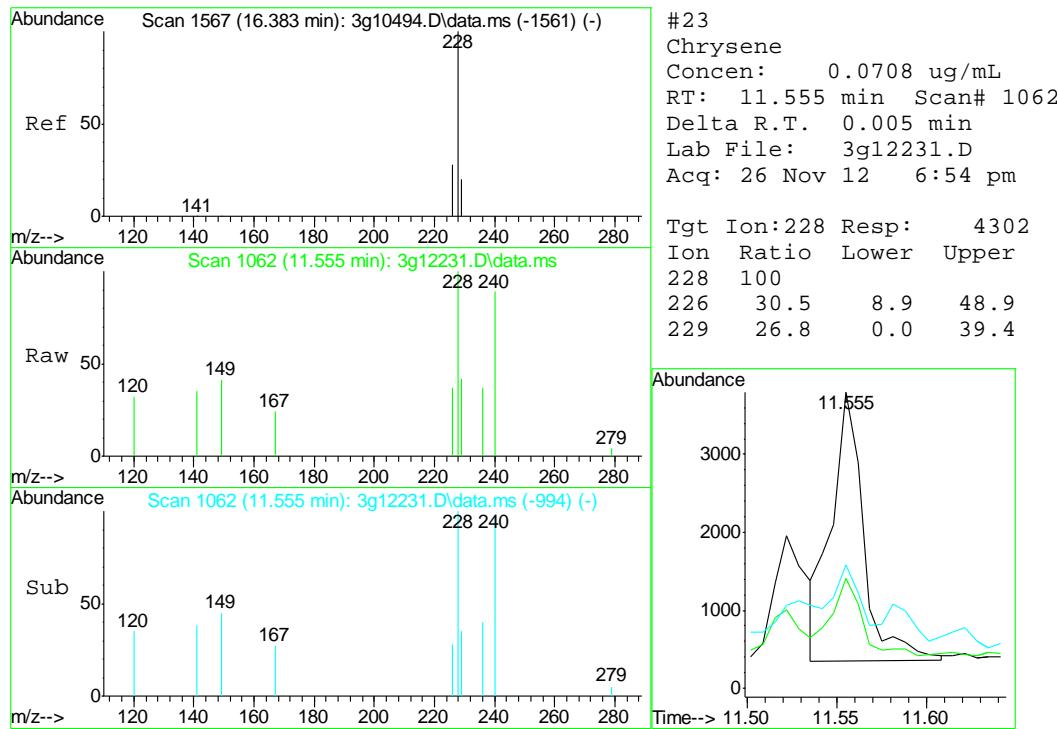


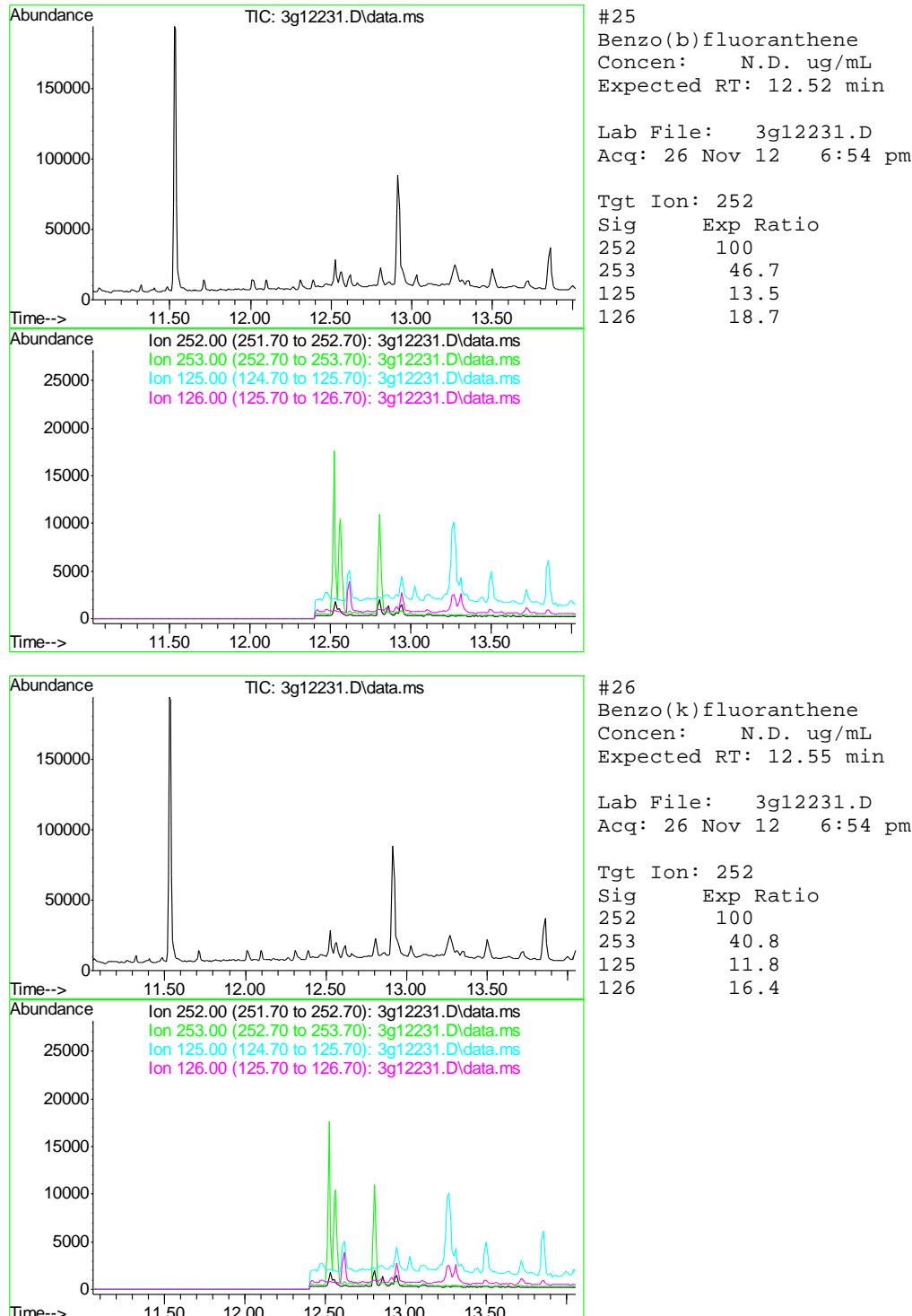


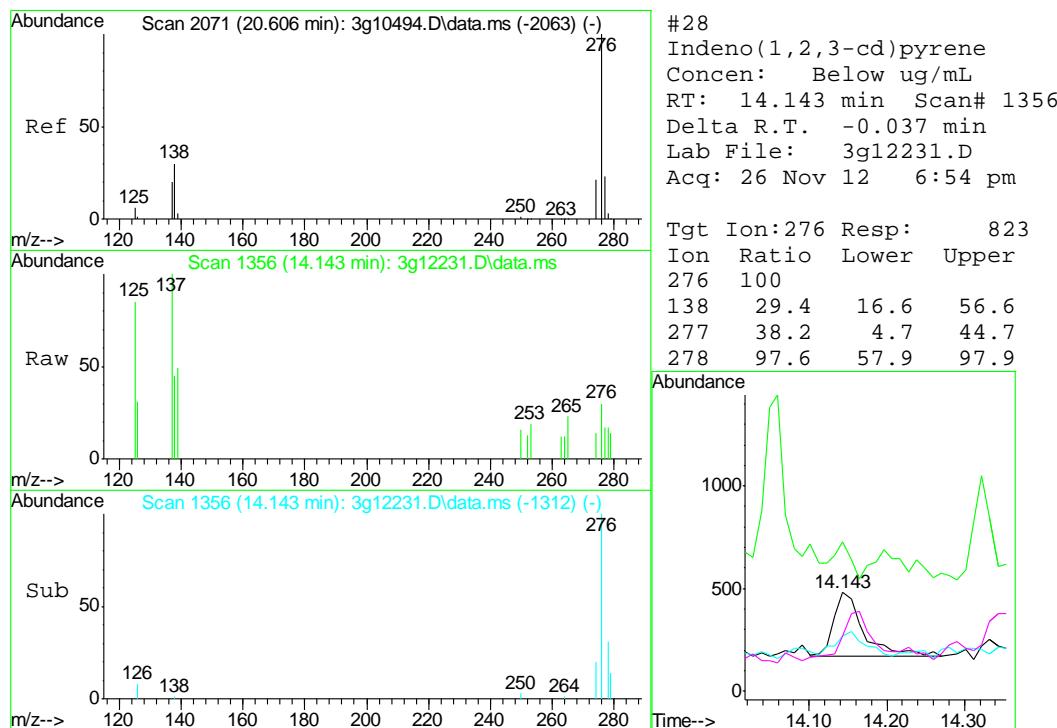
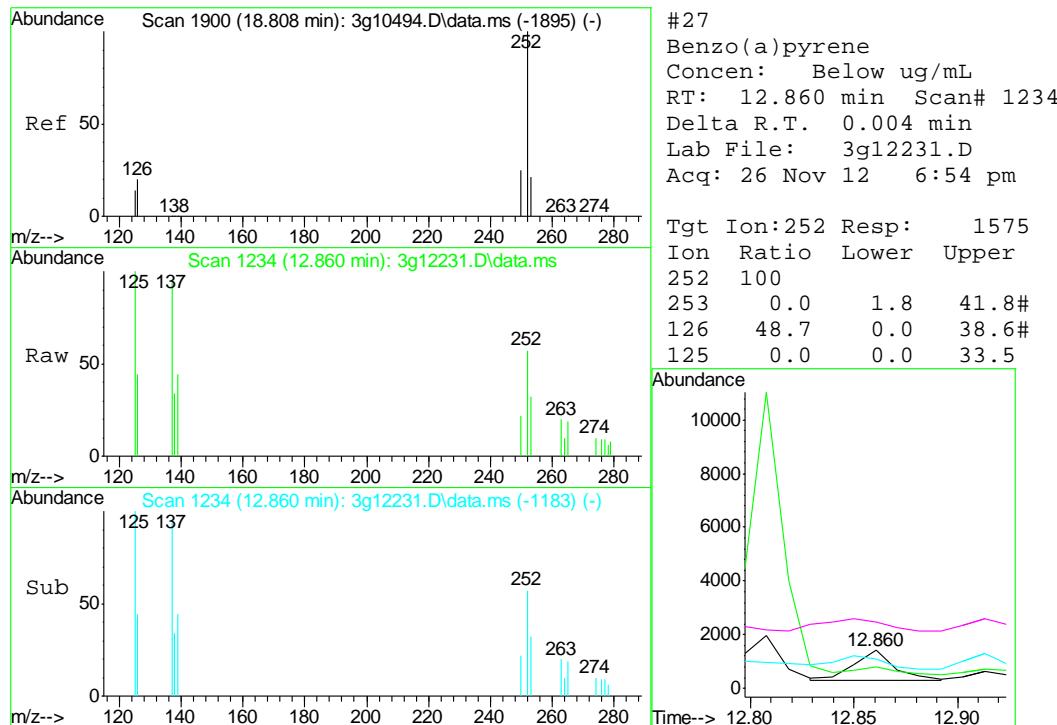


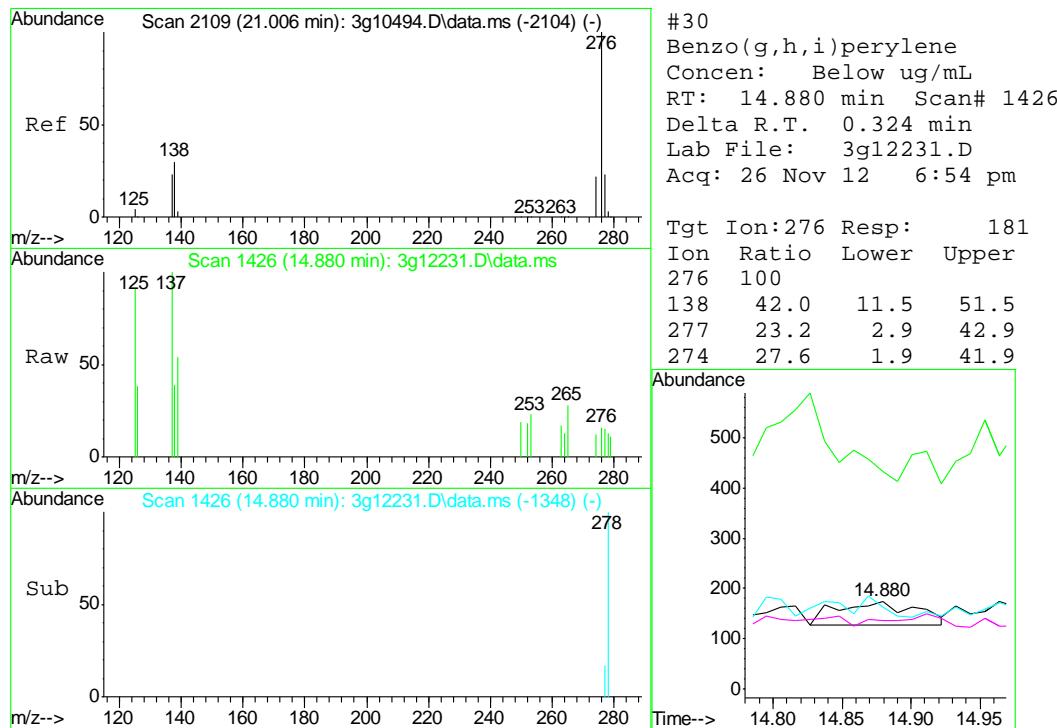
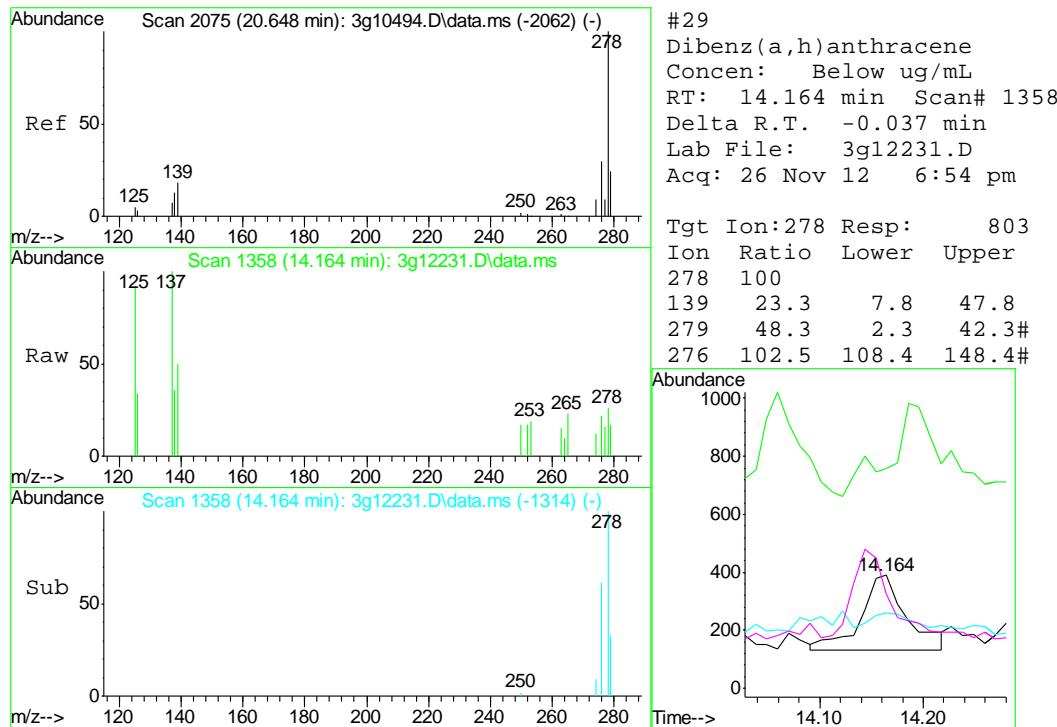












Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\112612\
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 Operator : SARAHM1
 Sample : OP6988-MB
 Misc : OP6988,E3G577,30.00,,,1,1
 ALS Vial : 7 Sample Multiplier: 1

Quant Time: Nov 27 09:22:34 2012
 Quant Method : C:\msdchem\1\METHODS\SIMPE3G574.M
 Quant Title : PAHSIM BASE
 QLast Update : Mon Nov 26 15:39:31 2012
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Naphthalene-d8	5.707	136	129576	4.0000	ug/mL	0.00
6) Acenaphthene-d10	7.412	164	82091	4.0000	ug/mL	0.00
15) Phenanthrene-d10	8.894	188	139846	4.0000	ug/mL	-0.06
19) Chrysene-d12	11.535	240	111025	4.0000	ug/mL	0.00
24) Perylene-d12	12.913	264	69568	4.0000	ug/mL	0.00

System Monitoring Compounds						
2) Nitrobenzene-d5	5.021	82	606301	48.6684	ug/mL	0.00
Spiked Amount	50.000	Range	25 - 135	Recovery	= 97.34%	
7) 2-Fluorobiphenyl	6.751	172	1487976	49.7336	ug/mL	0.00
Spiked Amount	50.000	Range	25 - 135	Recovery	= 99.46%	
21) Terphenyl-d14	10.493	244	659134	45.5508	ug/mL	0.00
Spiked Amount	50.000	Range	25 - 135	Recovery	= 91.10%	

Target Compounds					Qvalue
3) N-Nitrosodimethylamine	2.458	74	56	N.D.	
4) N-Nitrosodi-propylamine	0.000	70	0	N.D. d	
5) Naphthalene	5.719	128	385	N.D.	
8) 2-Methylnaphthalene	6.392	142	217	N.D.	
9) 1-Methylnaphthalene	6.492	142	148	N.D.	
10) Acenaphthylene	7.270	152	73	N.D.	
11) Acenaphthene	7.412	154	462	Below Cal #	70
12) Dibenzofuran	7.790	168	83	N.D.	
13) Fluorene	0.000	166	0	N.D. d	
14) Diphenylamine	0.000	169	0	N.D. d	
16) Phenanthrene	8.918	178	1004	Below Cal #	1
17) Anthracene	8.973	178	261	N.D.	
18) Fluoranthene	10.105	202	559	N.D.	
20) Pyrene	10.335	202	450	N.D.	
22) Benzo(a)anthracene	11.522	228	857	N.D.	
23) Chrysene	11.555	228	464	N.D.	
25) Benzo(b)fluoranthene	0.000	252	0	N.D. d	
26) Benzo(k)fluoranthene	12.534	252	410	N.D.	
27) Benzo(a)pyrene	12.860	252	352	N.D.	
28) Indeno(1,2,3-cd)pyrene	14.133	276	302	N.D.	
29) Dibenz(a,h)anthracene	14.154	278	241	N.D.	
30) Benzo(g,h,i)perylene	14.490	276	348	N.D.	

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

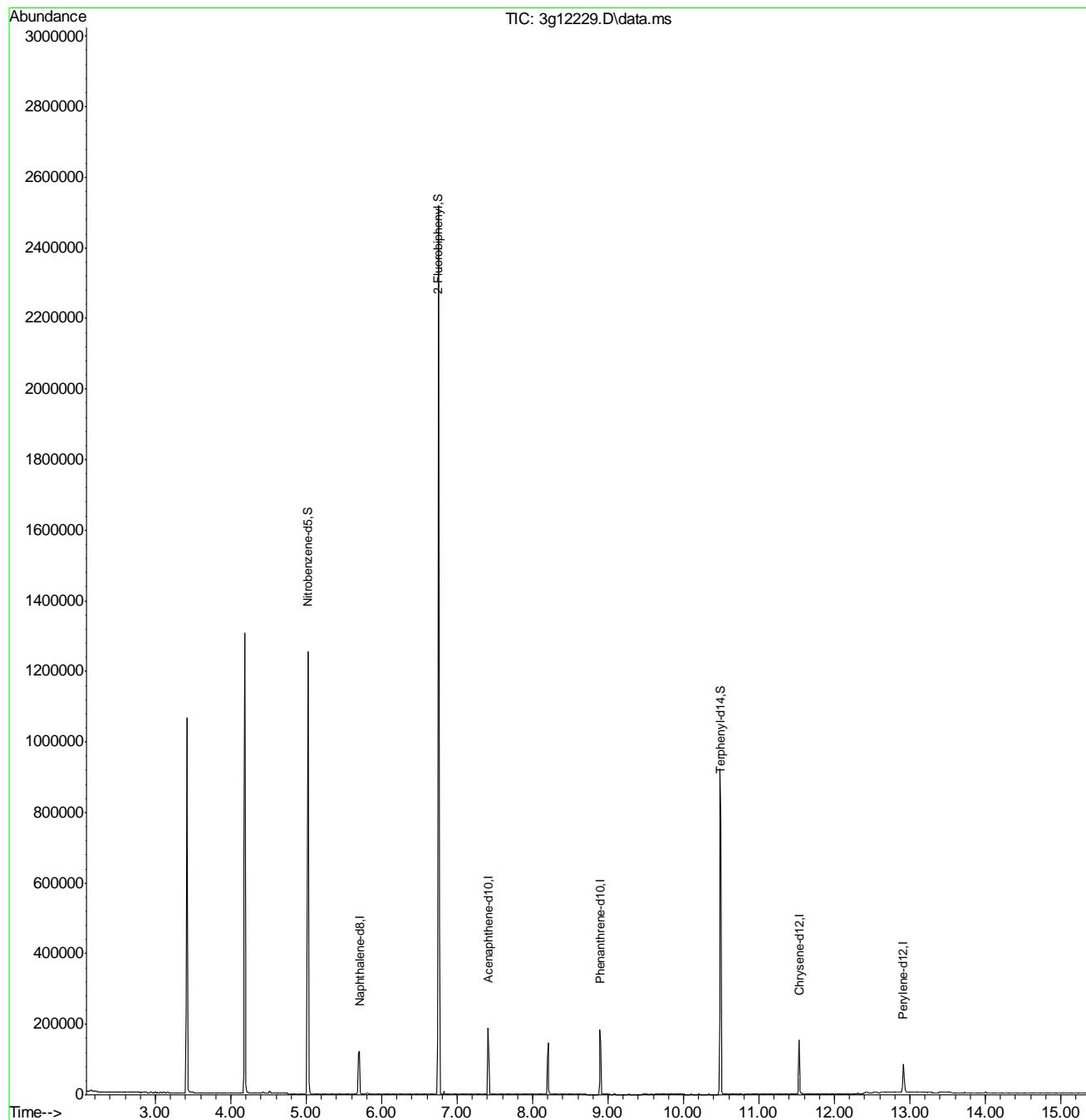
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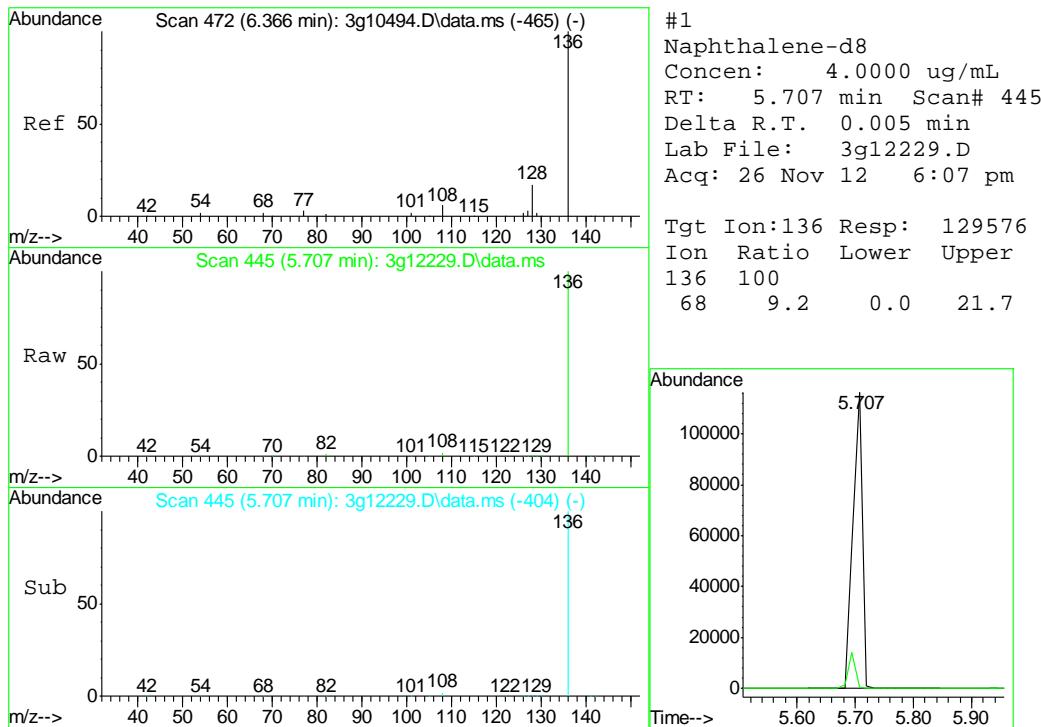
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Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\112612\
 Data File : 3g12229.D
 Acq On : 26 Nov 2012 6:07 pm
 Operator : SARAHM1
 Sample : OP6988-MB
 Misc : OP6988,E3G577,30.00,,,1,1
 ALS Vial : 7 Sample Multiplier: 1

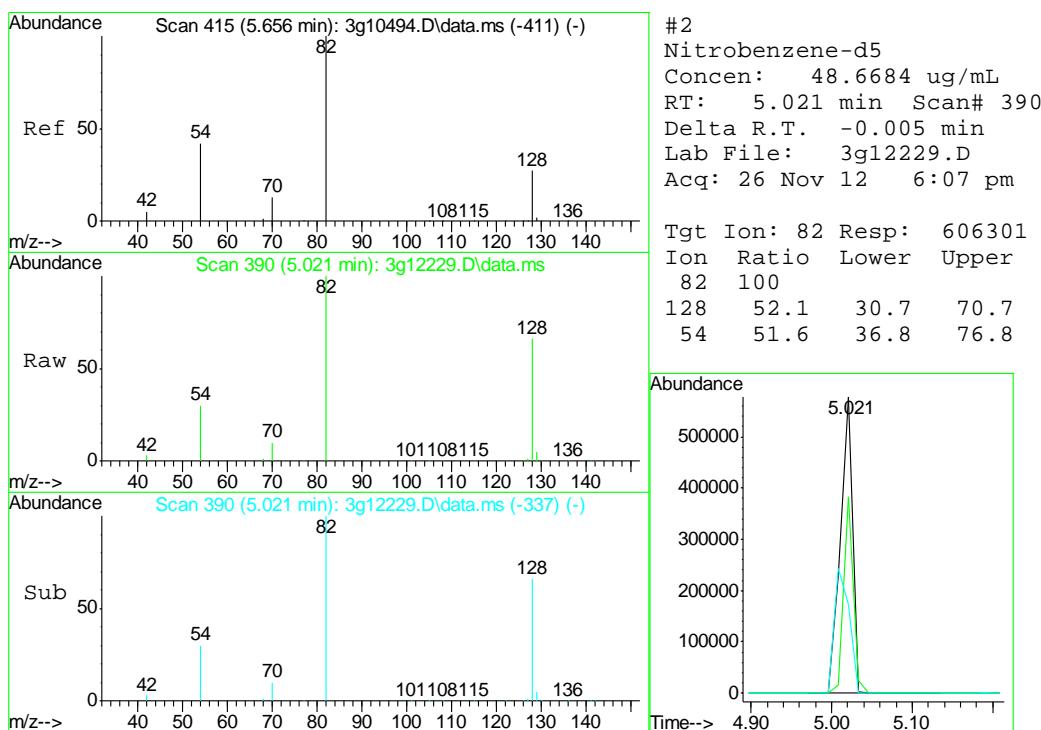
Quant Time: Nov 27 09:22:34 2012
 Quant Method : C:\msdchem\1\METHODS\SIMPE3G574.M
 Quant Title : PAHSIM BASE
 QLast Update : Mon Nov 26 15:39:31 2012
 Response via : Initial Calibration

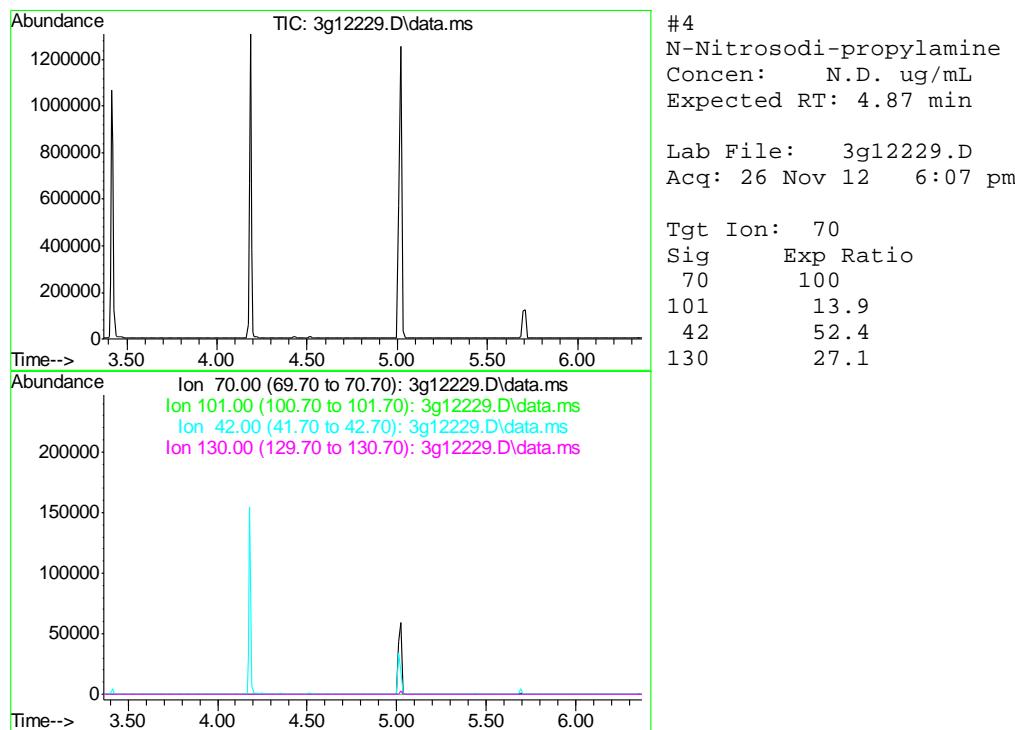
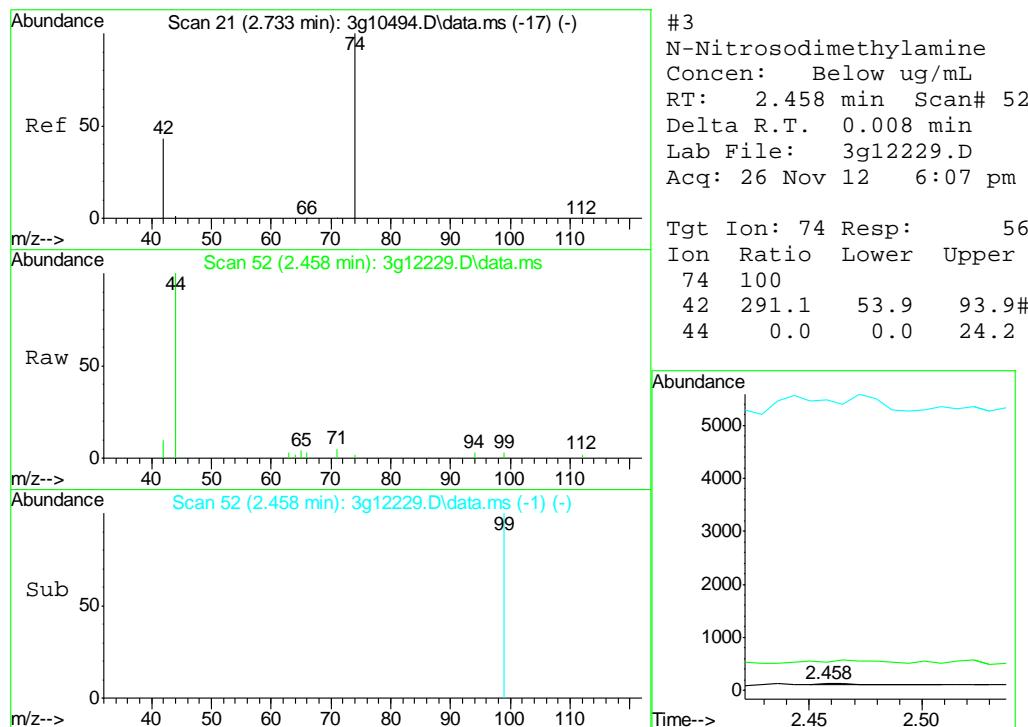


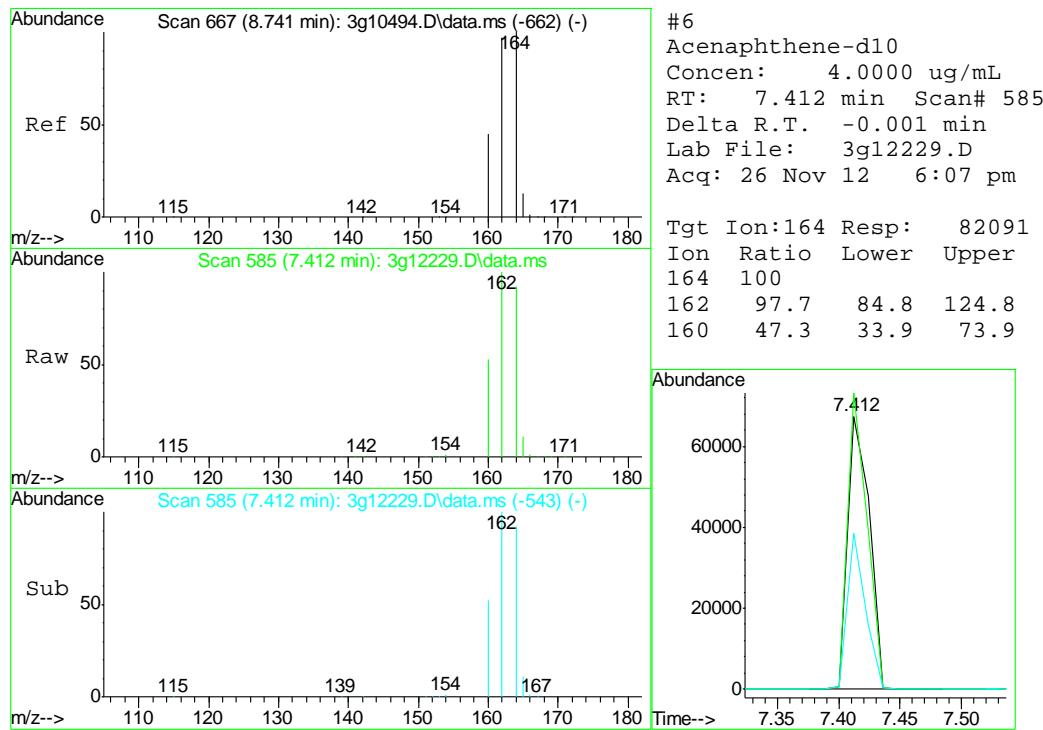
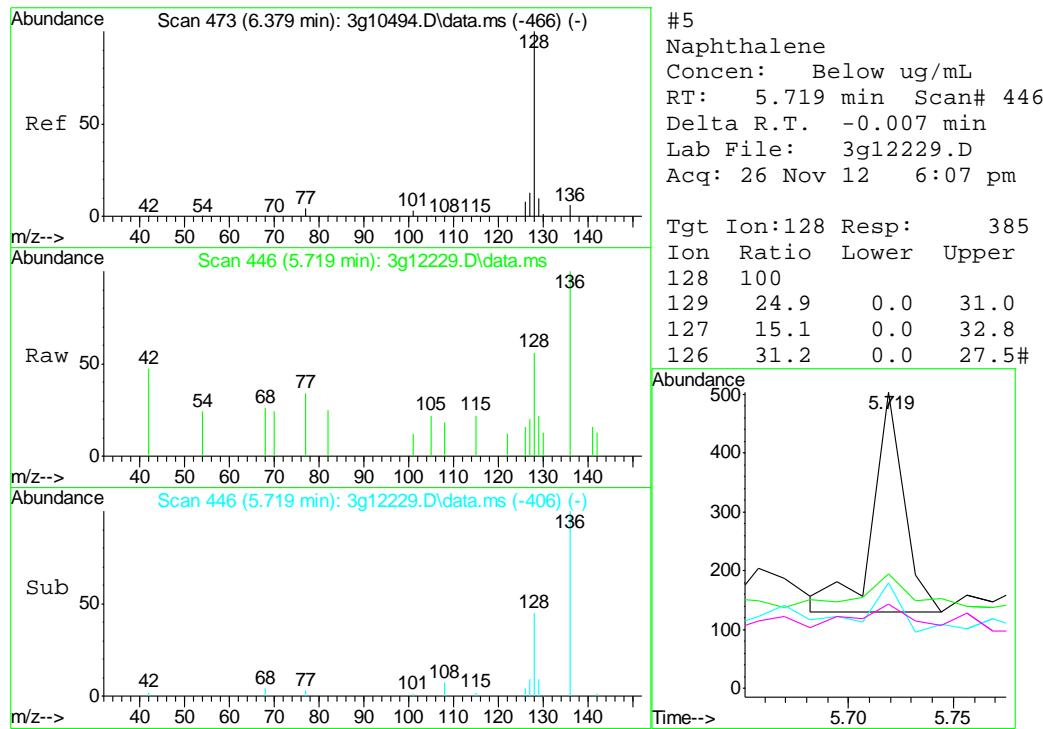


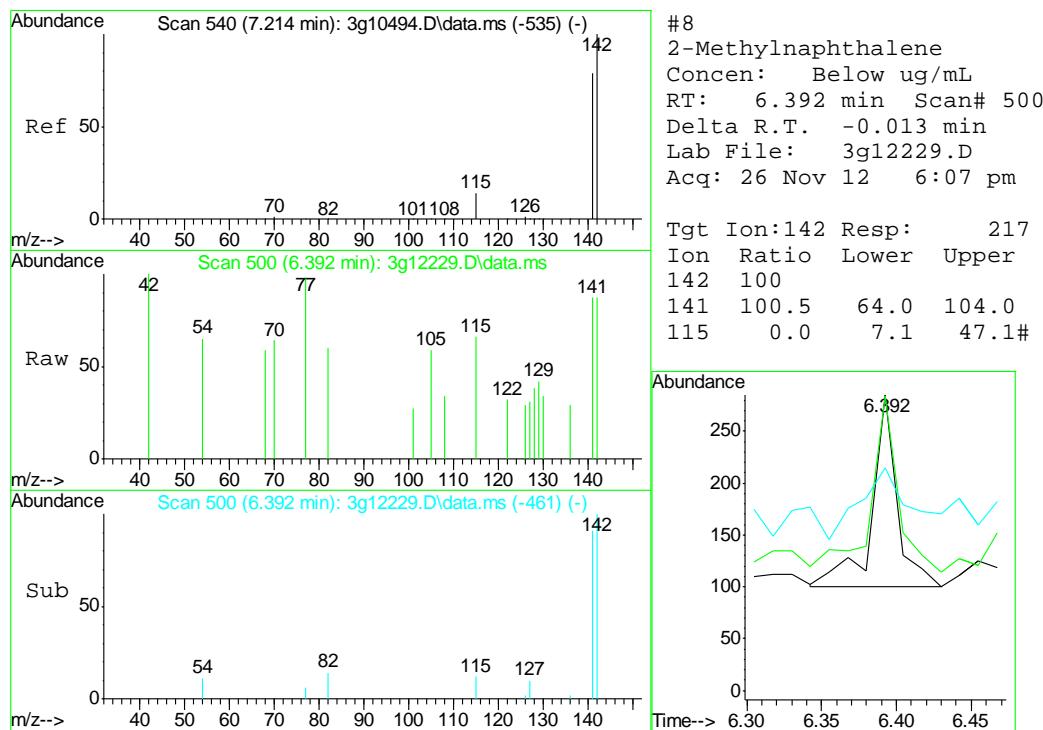
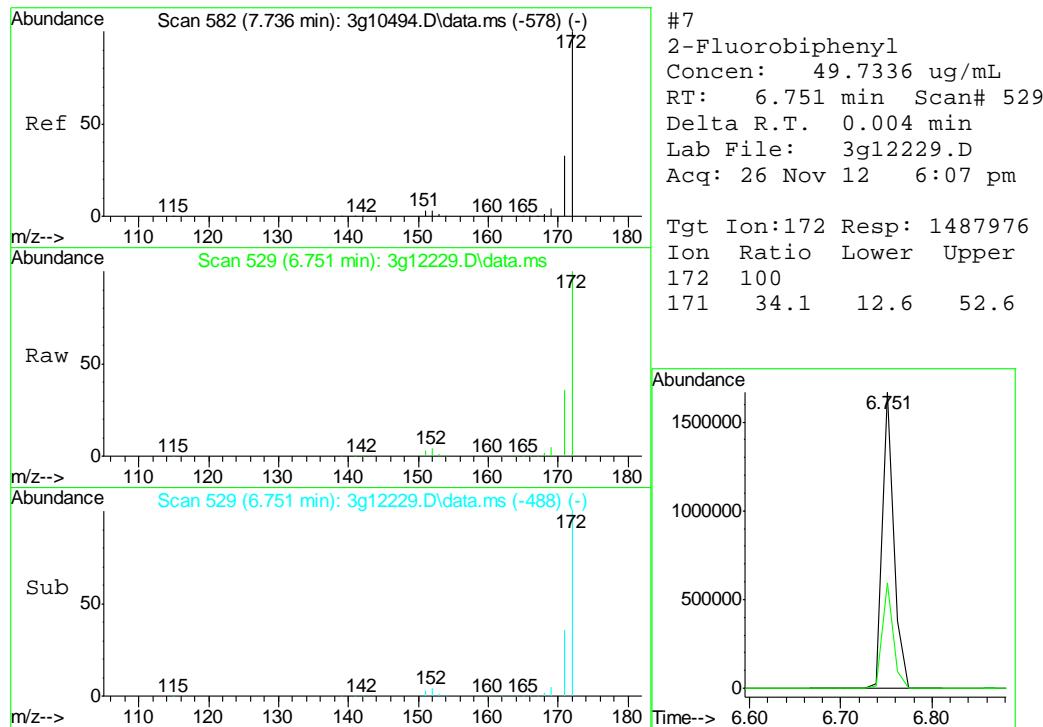
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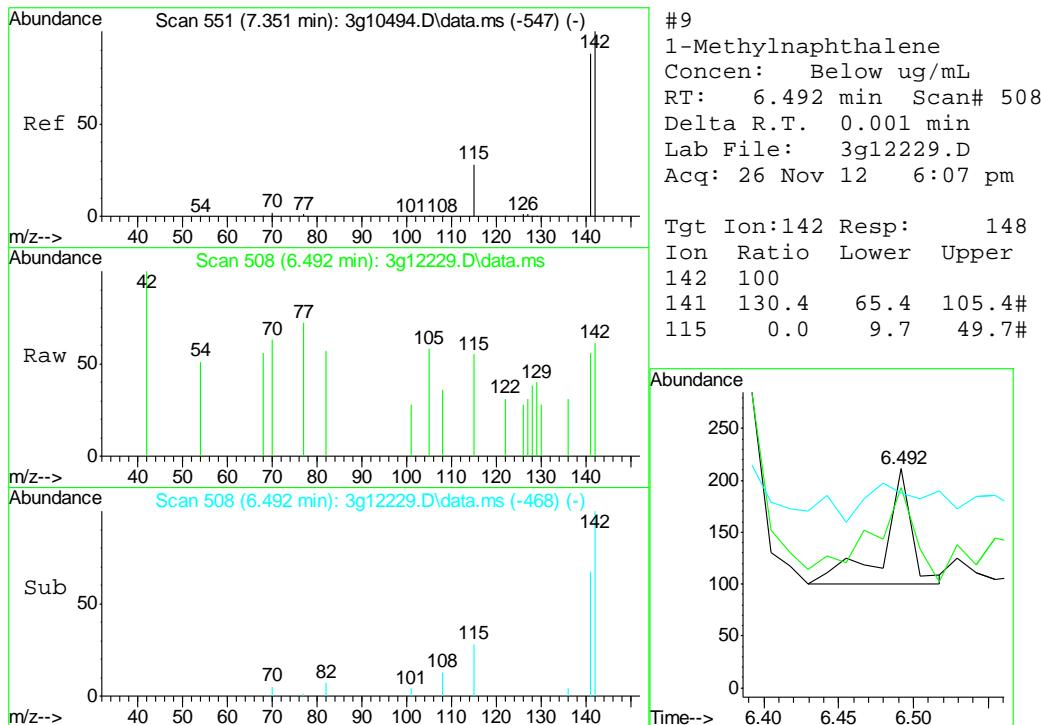
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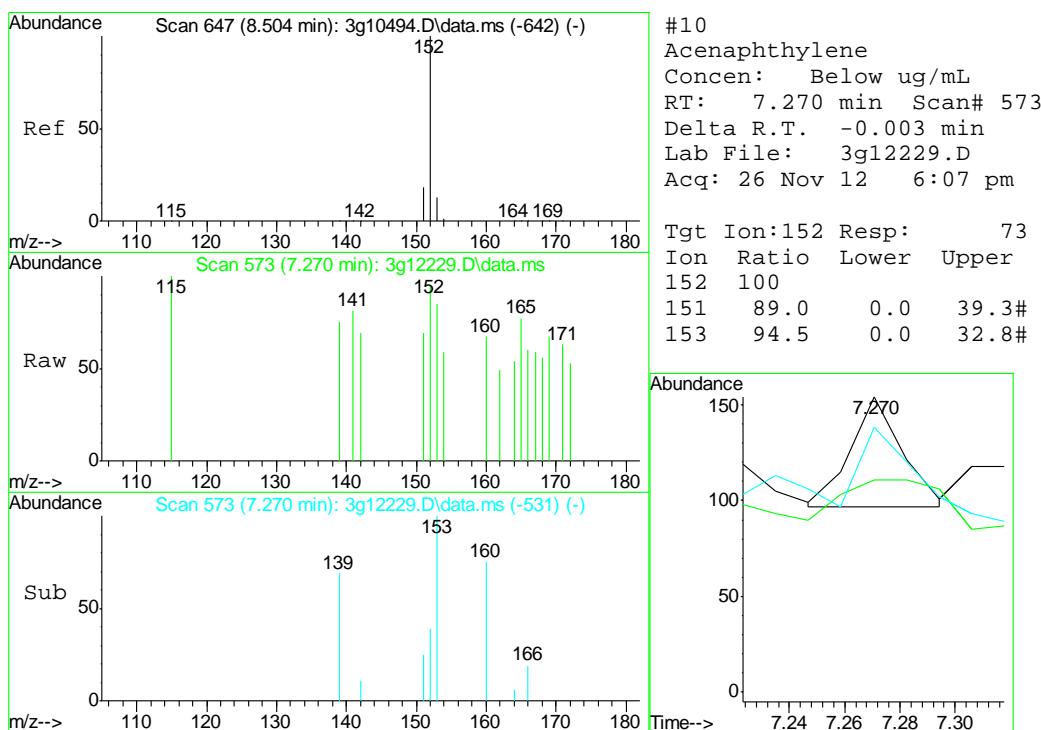


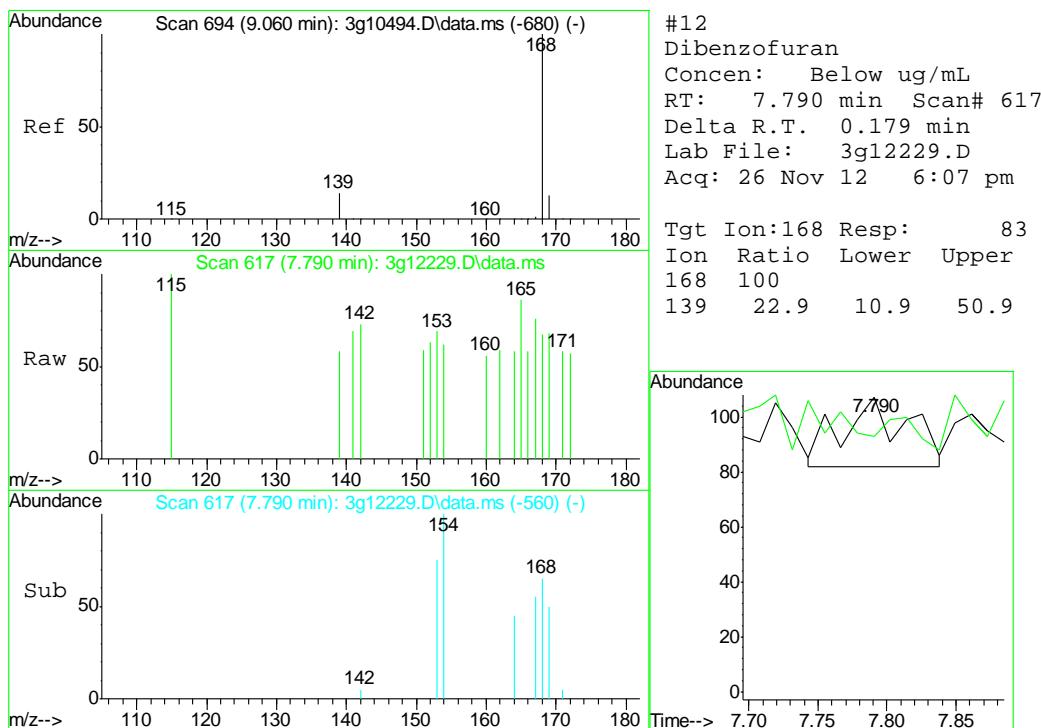
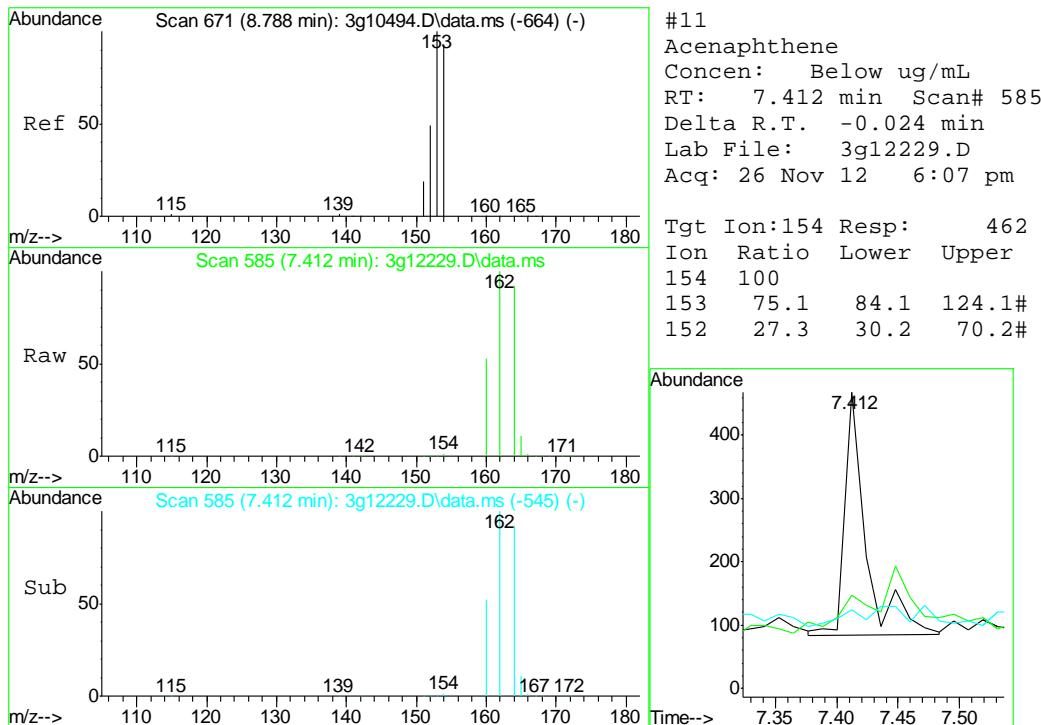


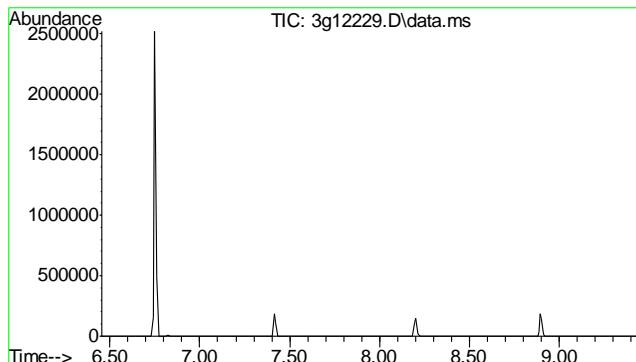


9.2.1

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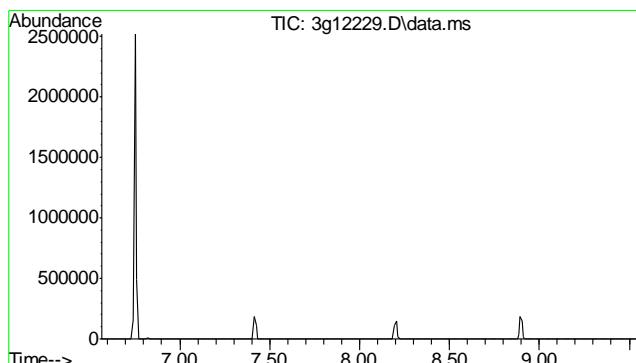
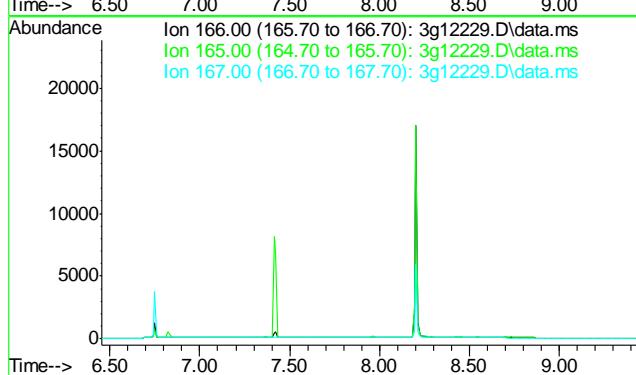




#13
Fluorene
Concen: N.D. ug/mL
Expected RT: 7.95 min

Lab File: 3g12229.D
Acq: 26 Nov 12 6:07 pm

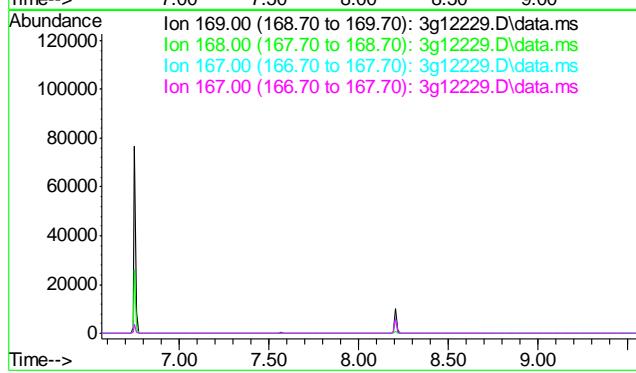
Tgt Ion: 166
Sig Exp Ratio
166 100
165 89.6
167 13.5

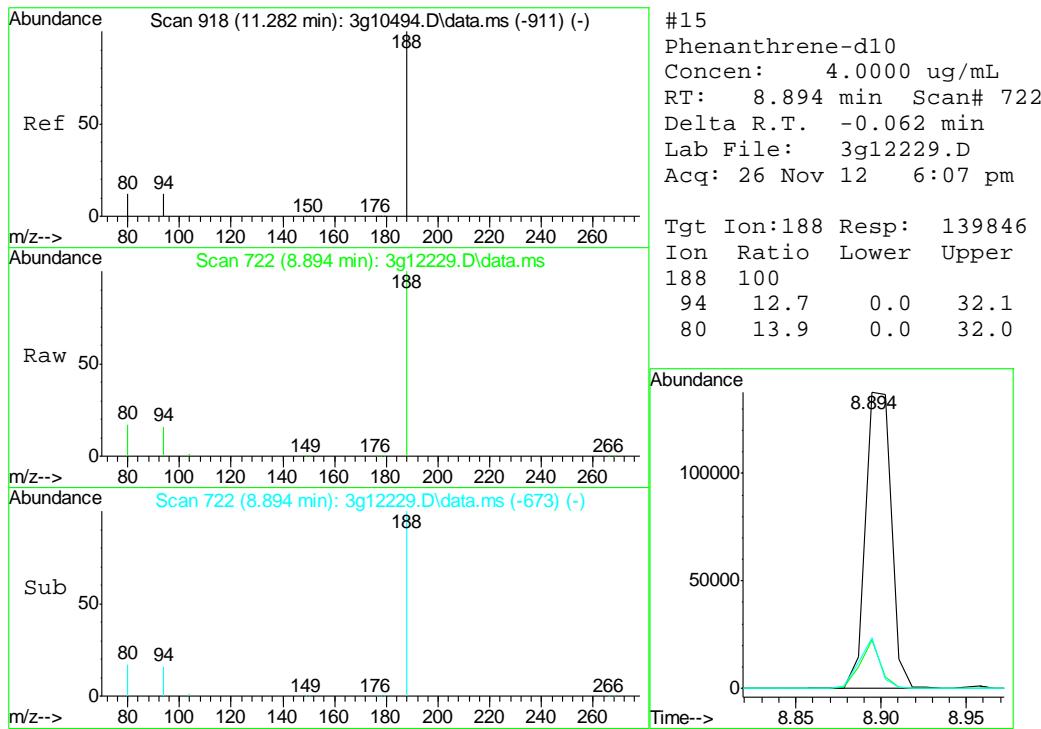


#14
Diphenylamine
Concen: N.D. ug/mL
Expected RT: 8.07 min

Lab File: 3g12229.D
Acq: 26 Nov 12 6:07 pm

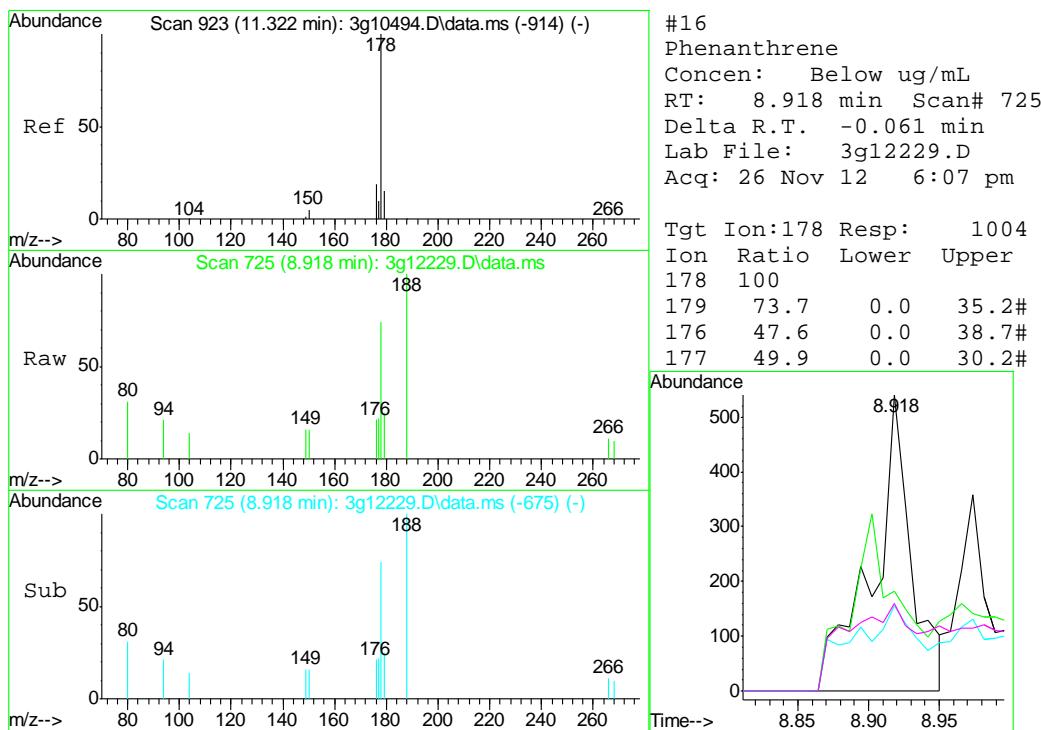
Tgt Ion: 169
Sig Exp Ratio
169 100
168 60.9
167 33.6
167 33.6

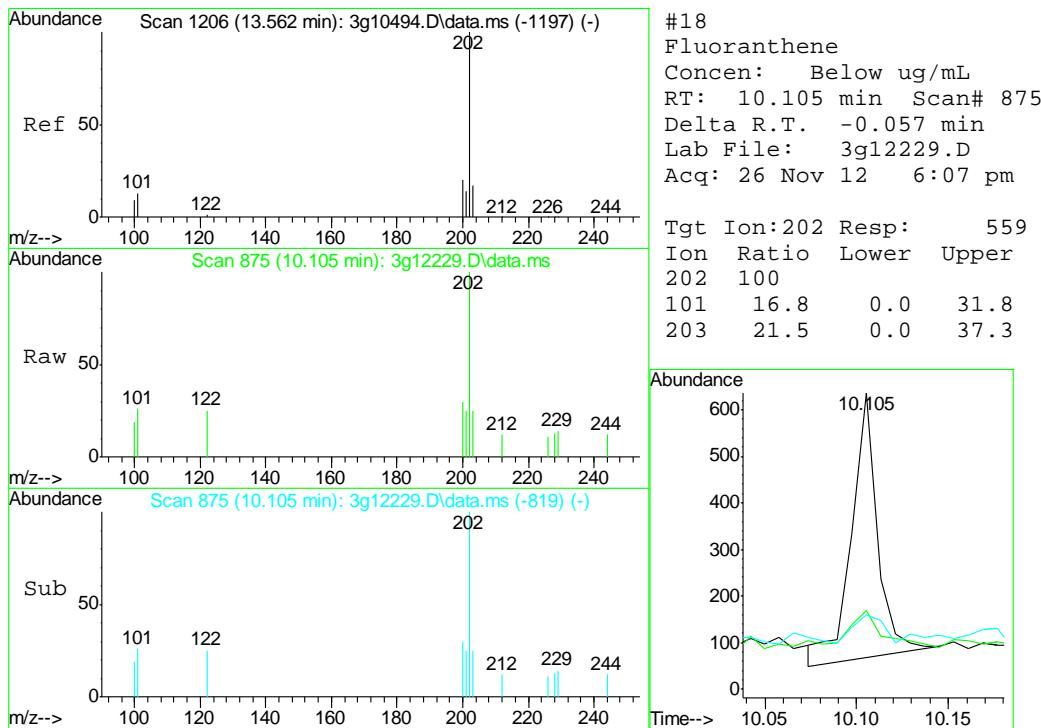
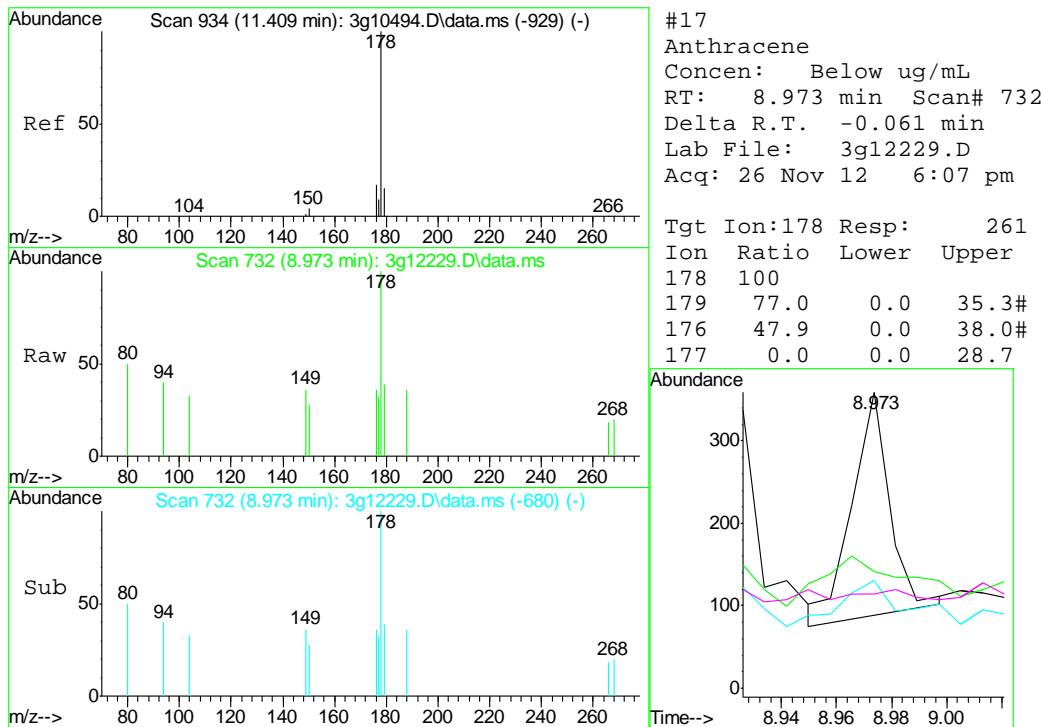


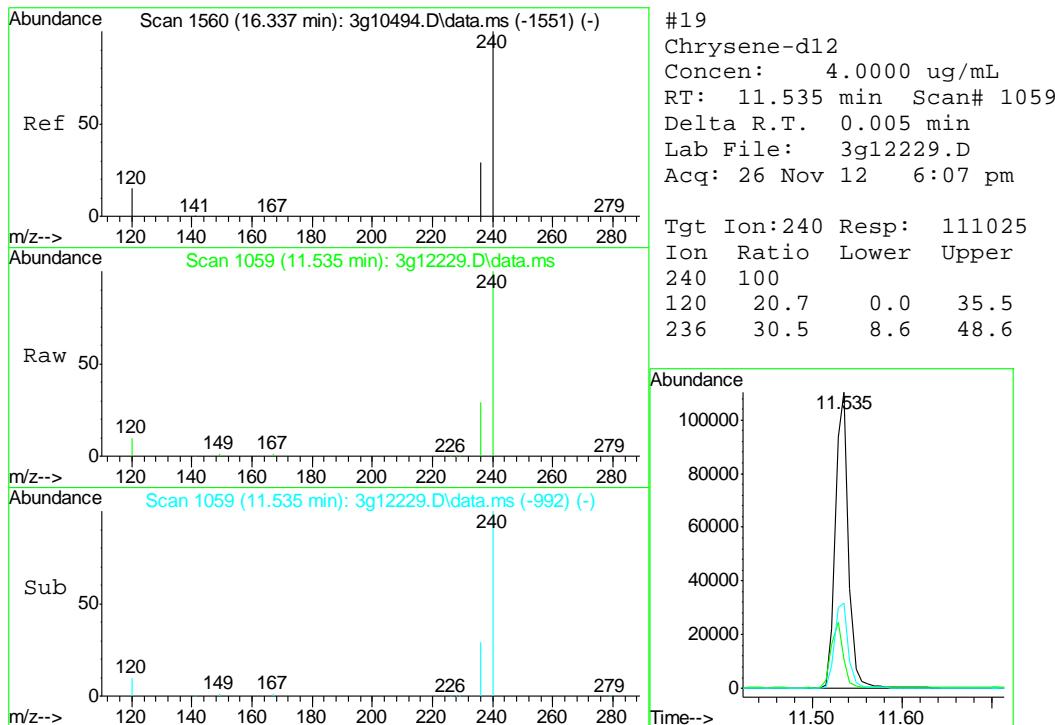


9.2.1

9

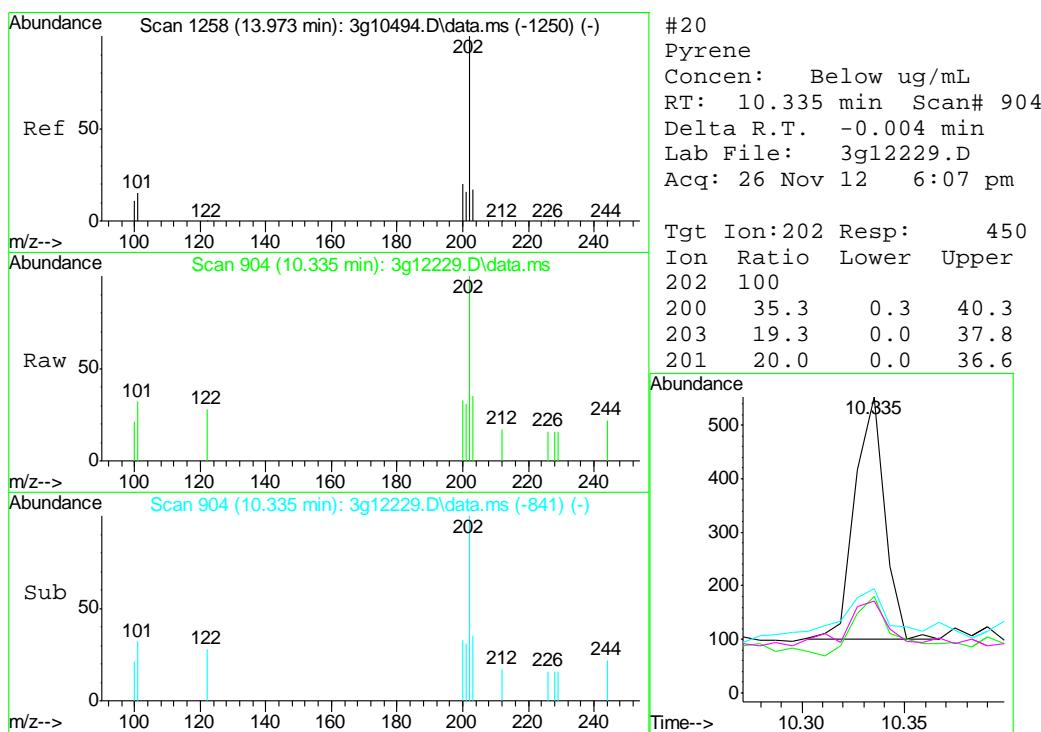


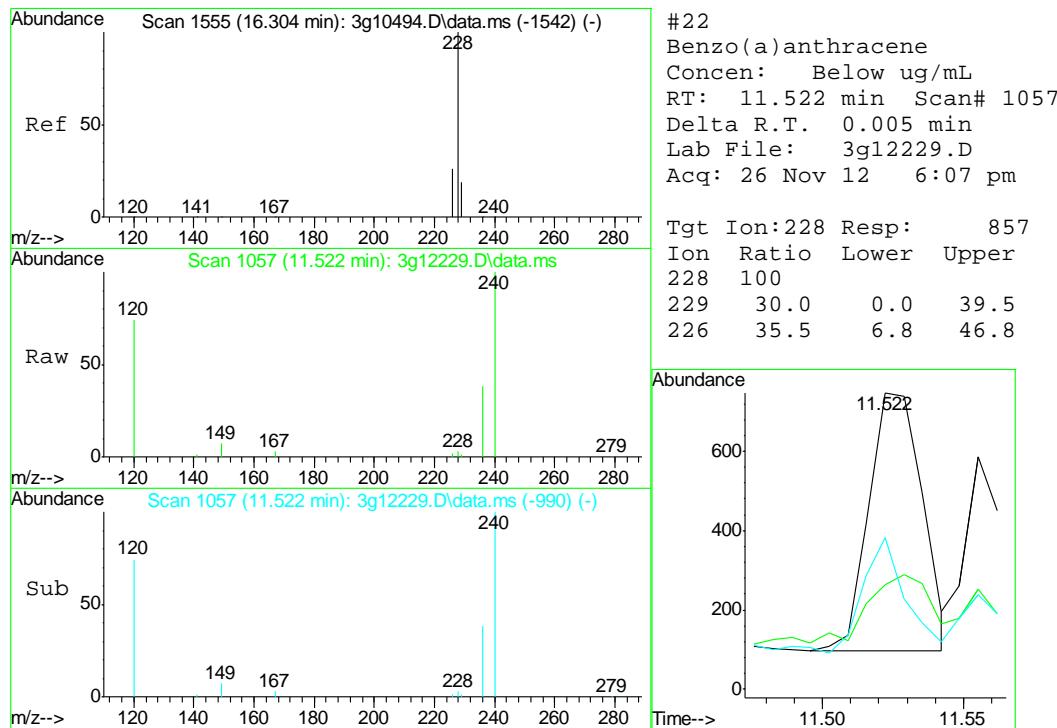
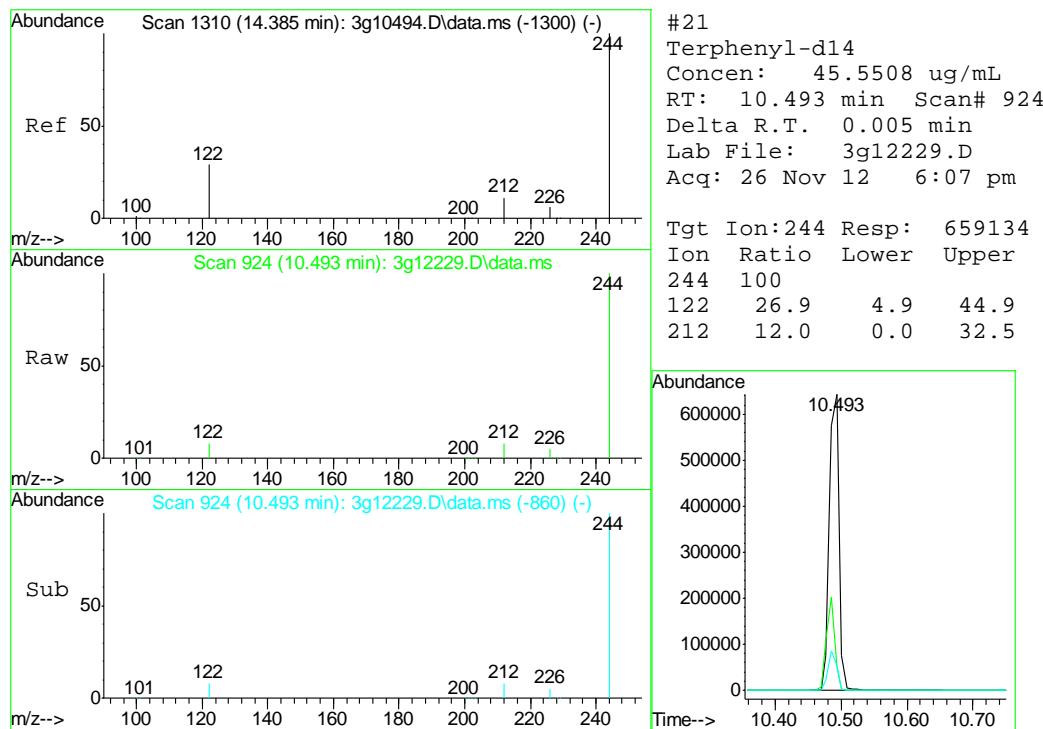


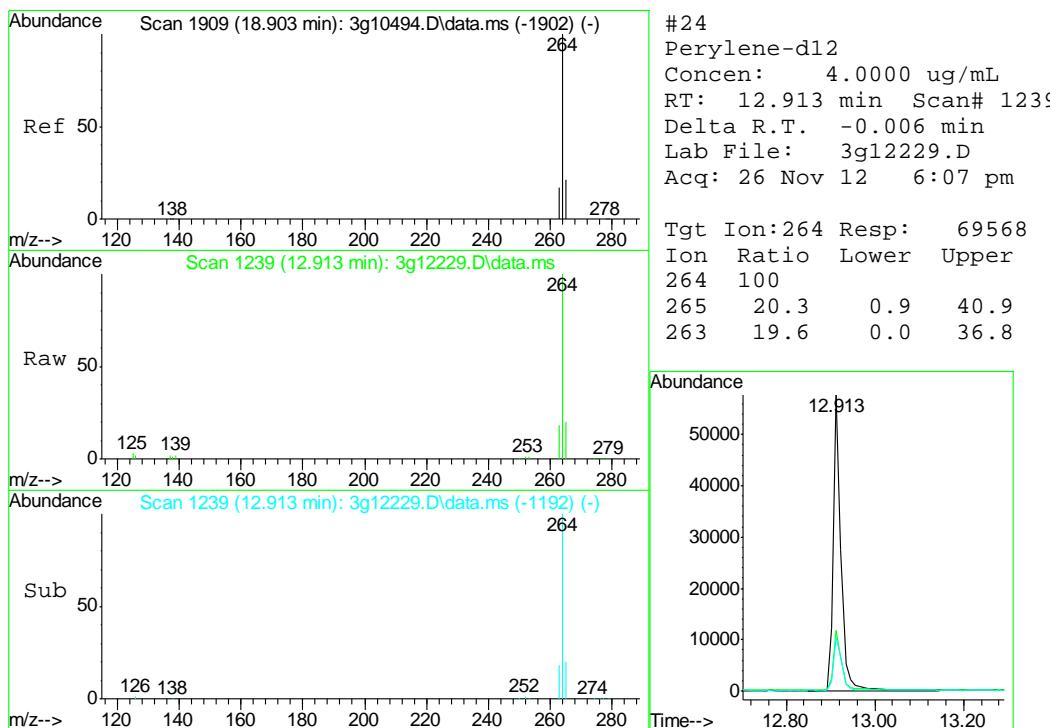
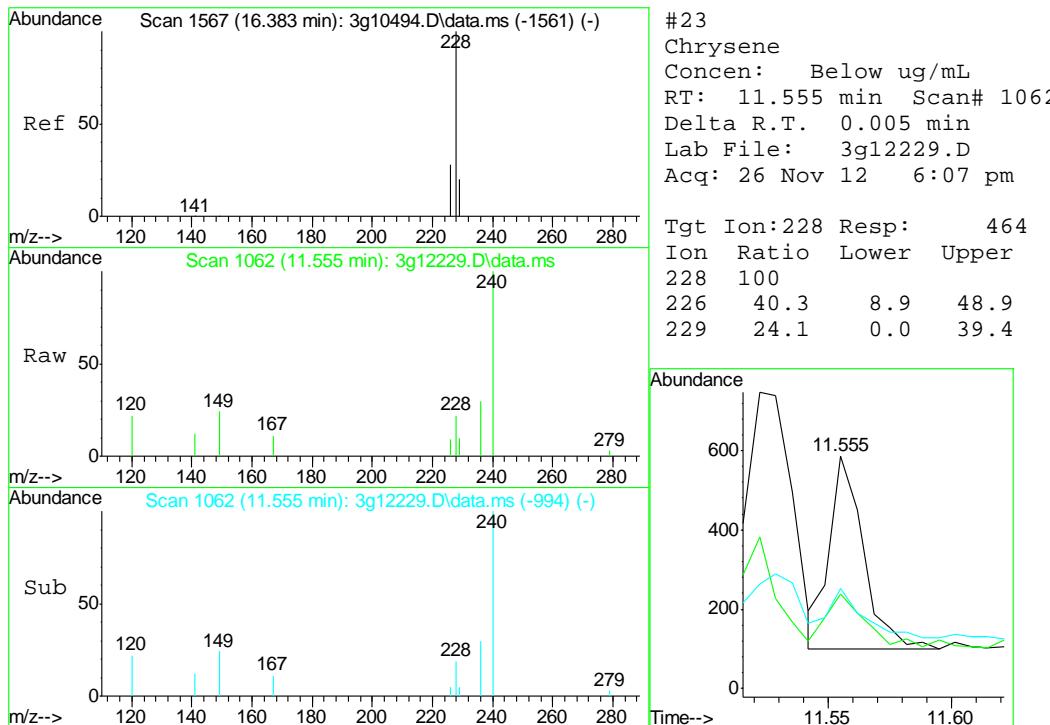


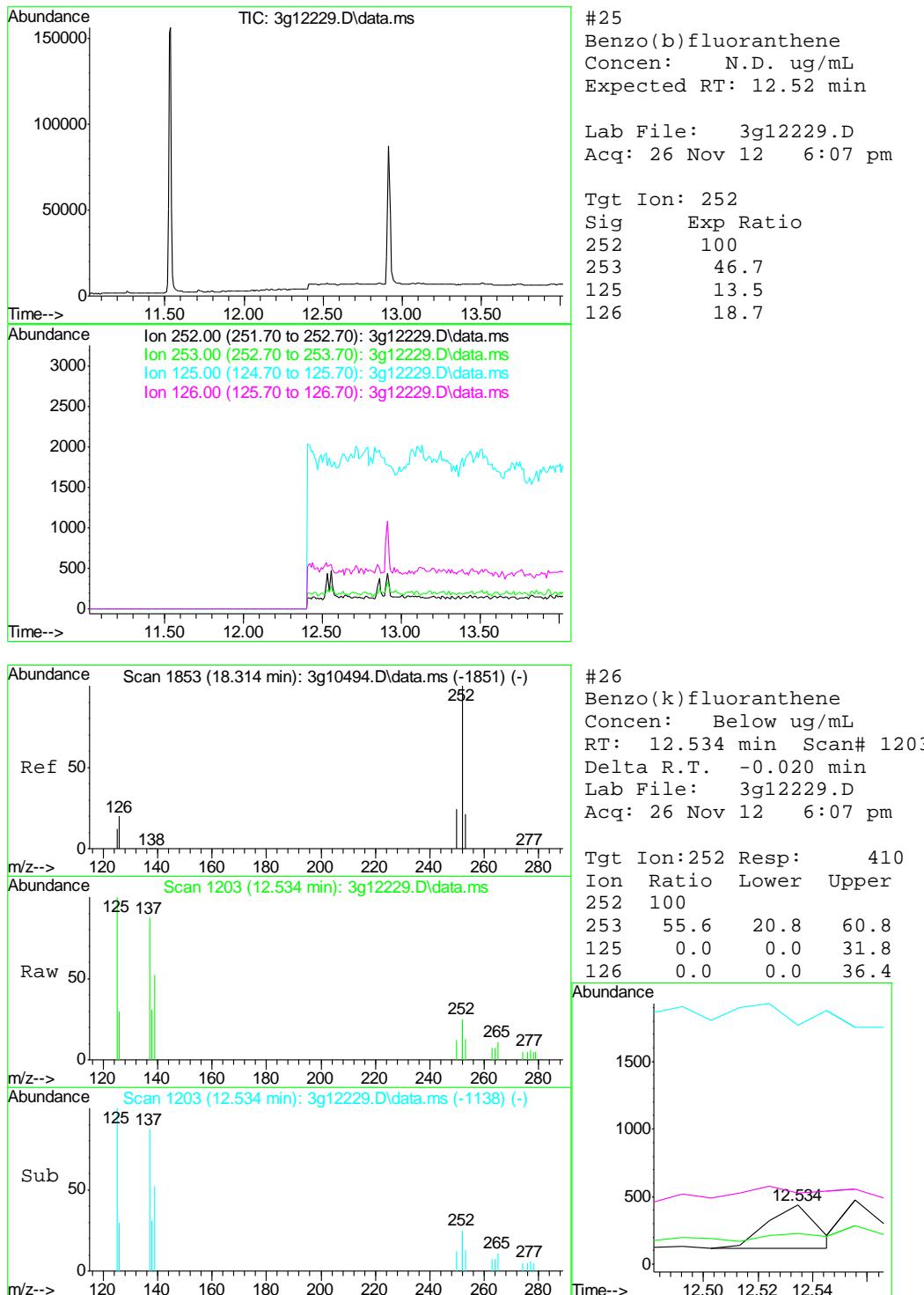
9.2.1

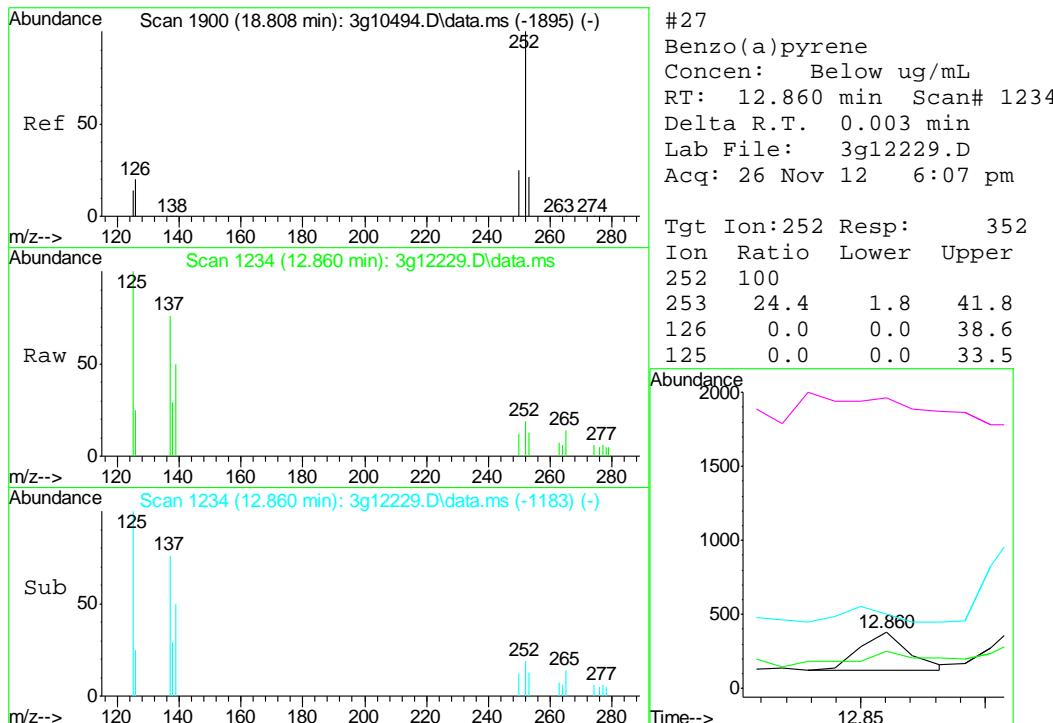
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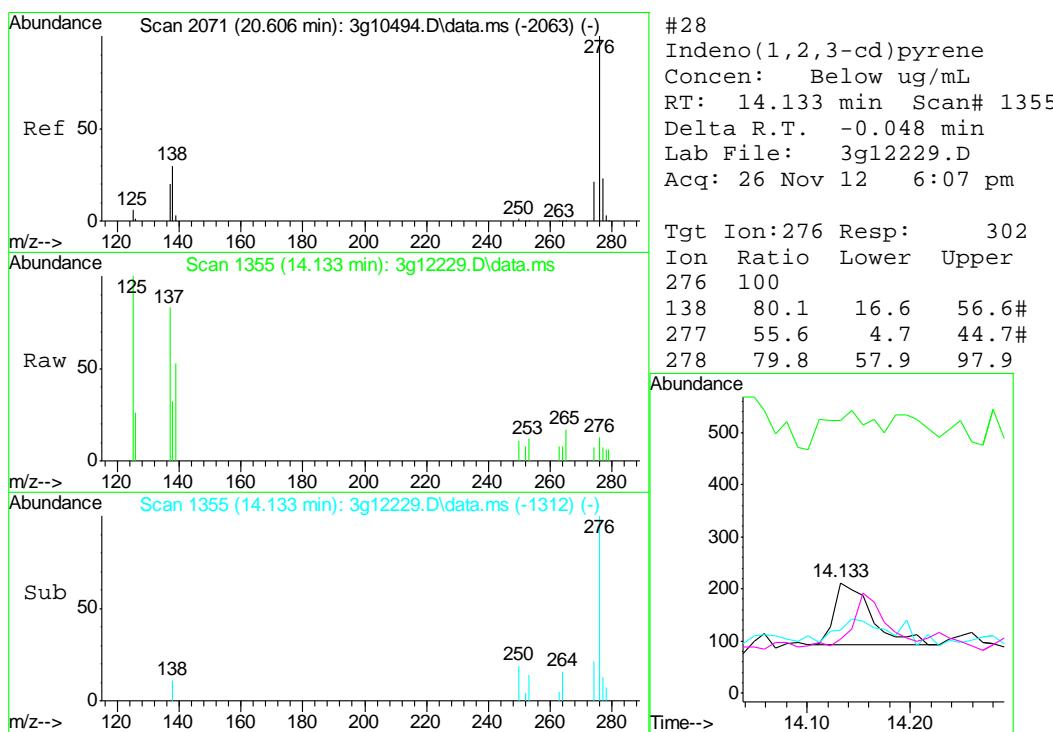


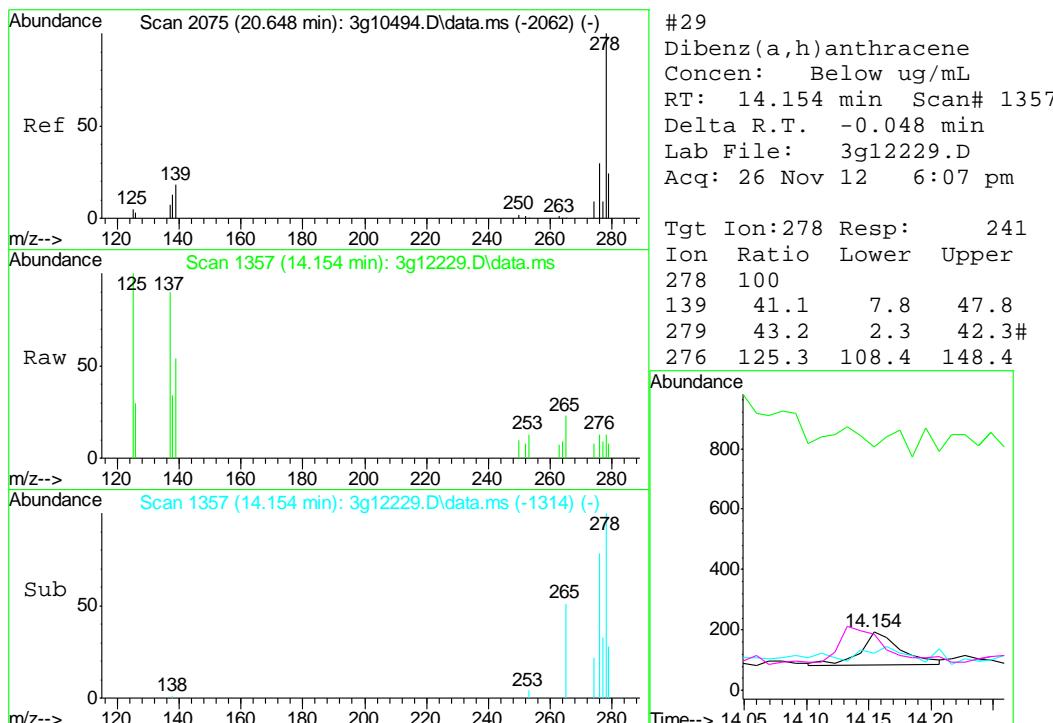




9.2.1

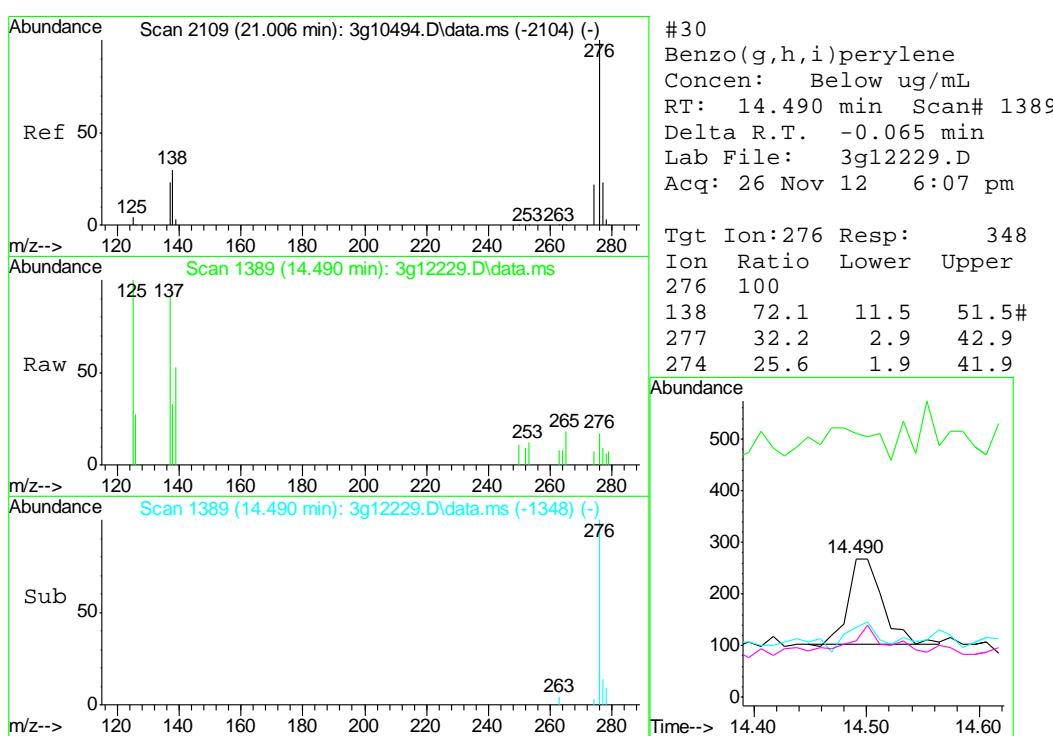
9





9.2.1

9





GC Volatiles

QC Data Summaries

Includes the following where applicable:

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

Method Blank Summary

Job Number: D41014

Account: XTOKWR XTO Energy

Project: PCU 296-5A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
GGB1010-MB	GB18499.D	1	11/16/12	SK	n/a	n/a	GGB1010

The QC reported here applies to the following samples:

Method: SW846 8015B

D41014-1

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	92% 60-140%

10.1.1

10

Blank Spike Summary

Page 1 of 1

Job Number: D41014

Account: XTOKWR XTO Energy

Project: PCU 296-5A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
GGB1010-BS	GB18500.D	1	11/16/12	SK	n/a	n/a	GGB1010

The QC reported here applies to the following samples:

Method: SW846 8015B

D41014-1

CAS No.	Compound	Spike mg/kg	BSP mg/kg	BSP %	Limits
	TPH-GRO (C6-C10)	110	124	113	70-130

CAS No.	Surrogate Recoveries	BSP	Limits
120-82-1	1,2,4-Trichlorobenzene	105%	60-140%

10.2.1
10

* = Outside of Control Limits.

Matrix Spike/Matrix Spike Duplicate Summary

Page 1 of 1

Job Number: D41014

Account: XTOKWR XTO Energy

Project: PCU 296-5A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
D41015-1MS	GB18502.D	1	11/16/12	SK	n/a	n/a	GGB1010
D41015-1MSD	GB18503.D	1	11/16/12	SK	n/a	n/a	GGB1010
D41015-1	GB18501.D	1	11/16/12	SK	n/a	n/a	GGB1010

The QC reported here applies to the following samples:

Method: SW846 8015B

D41014-1

CAS No.	Compound	D41015-1		Spike	MS	MS	MSD	MSD	RPD	Limits Rec/RPD
		mg/kg	Q	mg/kg	mg/kg	%	mg/kg	%		
	TPH-GRO (C6-C10)	ND		128	146	114	146	114	0	70-130/30

CAS No.	Surrogate Recoveries	MS	MSD	D41015-1	Limits
120-82-1	1,2,4-Trichlorobenzene	109%	111%	93%	60-140%

* = Outside of Control Limits.

10.3.1
10



GC Volatiles

Raw Data

Quantitation Report (QT Reviewed)

Signal #1 : Y:\1\DATA\111612\GB18507.D\FID1A.CH Vial: 11
 Signal #2 : Y:\1\DATA\111612\GB18507.D\FID2B.CH
 Acq On : 16 Nov 2012 10:34 pm Operator: StephK
 Sample : D41014-1, 50X Inst : GC/MS Ins
 Misc : GC3242,GGB1010,5.068,,100,5,1 Multiplr: 1.00
 IntFile Signal #1: TVH1.E IntFile Signal #2: FB2.E
 Quant Time: Nov 17 09:39:26 2012 Quant Results File: TB868GB868SOIL.RES

Quant Method : C:\MSDCHEM\1...\TB868GB868SOIL.M (Chemstation Integrator)
 Title : 8015B/8021B TVH/BTEX
 Last Update : Fri Nov 16 18:14:55 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	14.38	2867749	91.522 %
10) S	1,2,4-Trichlorobenzene (P)	14.38	15266238	93.930 %

Target Compounds

1) H	TVH-Gasoline	7.23	4368751	<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	177320	0.447 ug/L
7) T	Ethylbenzene	0.00	0	N.D. ug/L d
8) T	m,p-Xylene	10.48	174040	0.103 ug/L
9) T	o-Xylene	0.00	0	N.D. ug/L d
11) T	Naphthalene	14.56	233105	1.181 ug/L

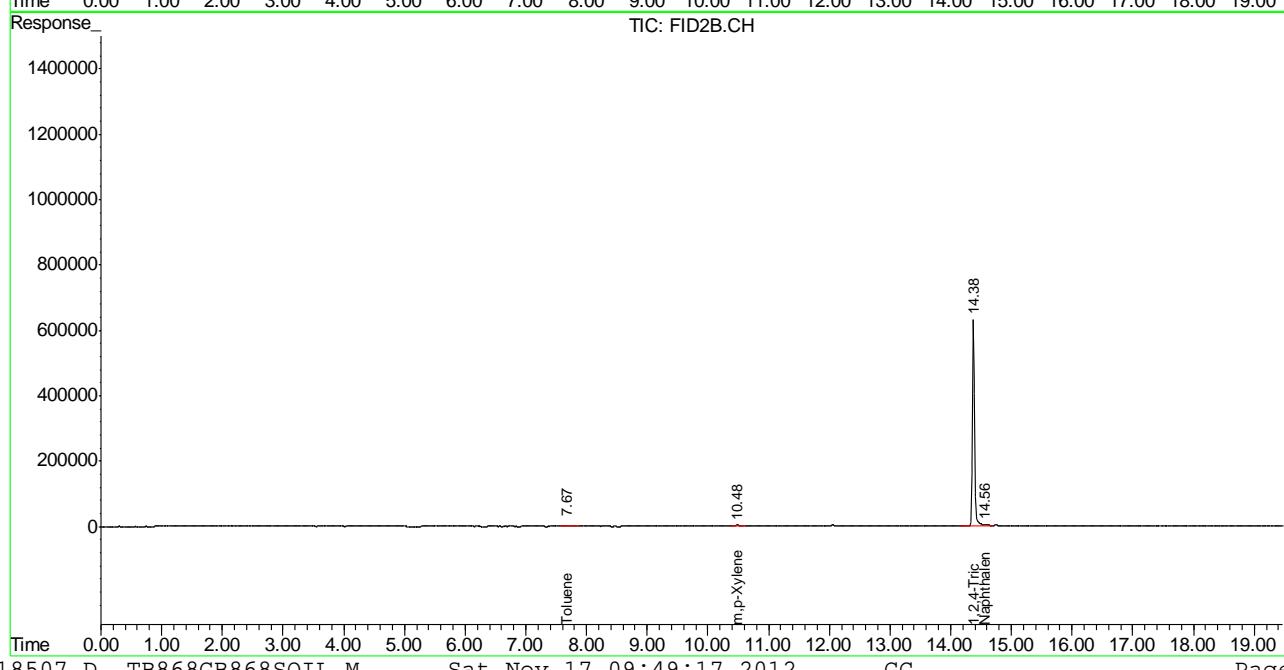
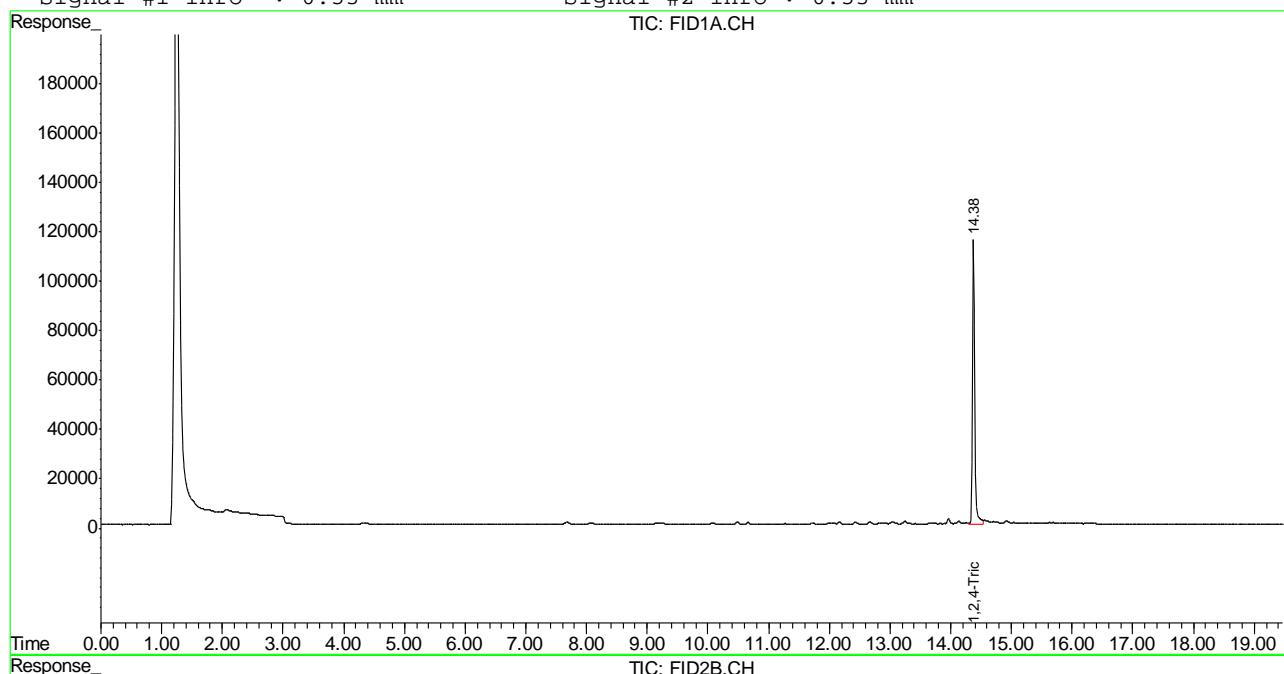
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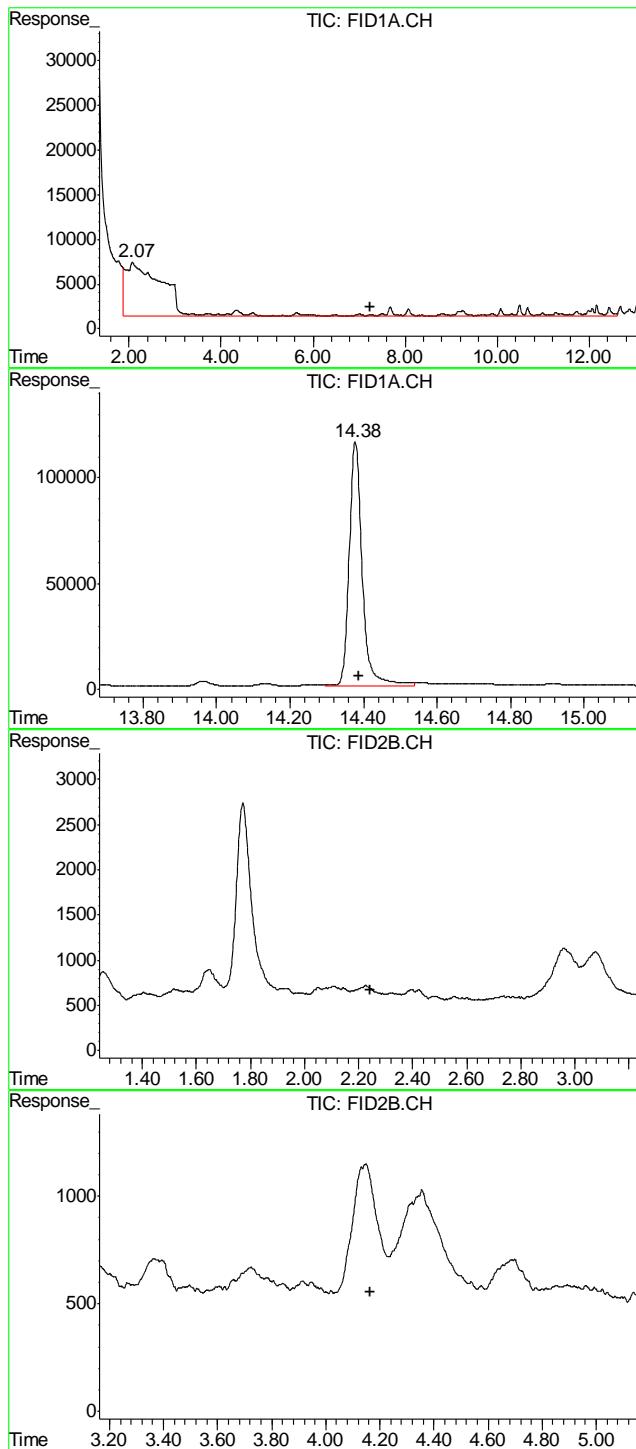
Quantitation Report (QT Reviewed)

Signal #1 : Y:\1\DATA\111612\GB18507.D\FID1A.CH Vial: 11
 Signal #2 : Y:\1\DATA\111612\GB18507.D\FID2B.CH
 Acq On : 16 Nov 2012 10:34 pm Operator: StephK
 Sample : D41014-1, 50X Inst : GC/MS Ins
 Misc : GC3242,GGB1010,,5.068,,100,5,1 Multiplr: 1.00
 IntFile Signal #1: TVH1.E IntFile Signal #2: FB2.E
 Quant Time: Nov 17 9:43 2012 Quant Results File: TB868GB868SOIL.RES

Quant Method : C:\MSDCHEM\1...\TB868GB868SOIL.M (Chemstation Integrator)
 Title : 8015B/8021B TVH/BTEX
 Last Update : Fri Nov 16 18:14:55 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: 4368751
 Conc: N.D.

#2 1,2,4-Trichlorobenzene

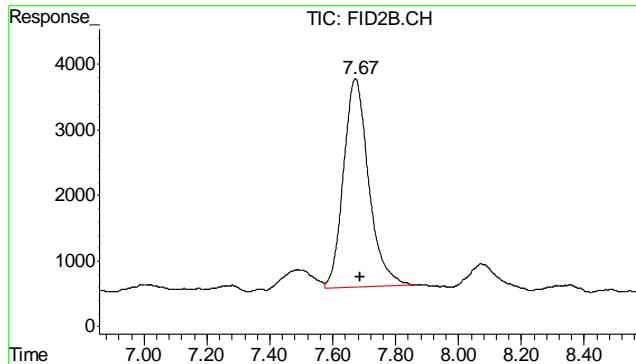
R.T.: 14.378 min
 Delta R.T.: -0.010 min
 Response: 2867749
 Conc: 91.52 %

#4 Methyl-t-butyl-ether

R.T.: 0.000 min
 Exp R.T. : 2.242 min
 Response: 0
 Conc: N.D.

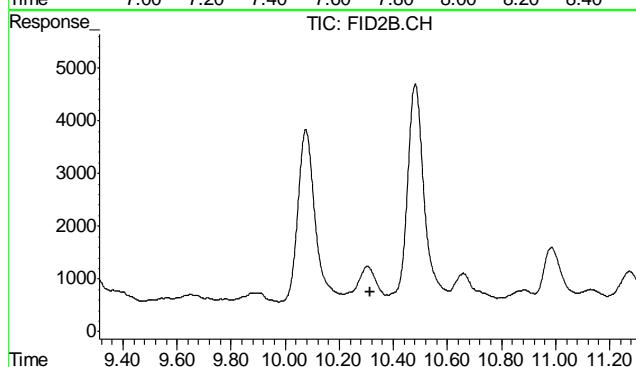
#5 Benzene

R.T.: 0.000 min
 Exp R.T. : 4.163 min
 Response: 0
 Conc: N.D.



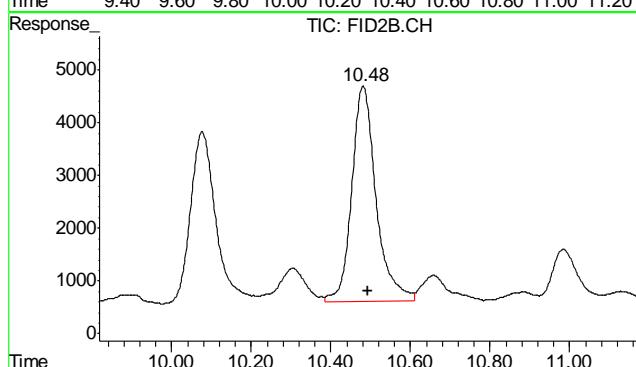
#6 Toluene

R.T.: 7.672 min
Delta R.T.: -0.016 min
Response: 177320
Conc: 0.45 ug/L



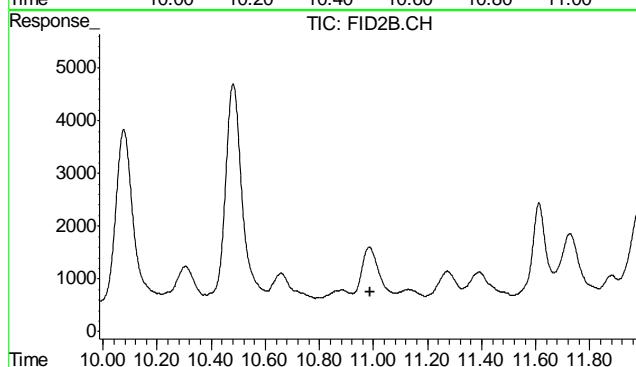
#7 Ethylbenzene

R.T.: 0.000 min
Exp R.T.: 10.313 min
Response: 0
Conc: N.D.



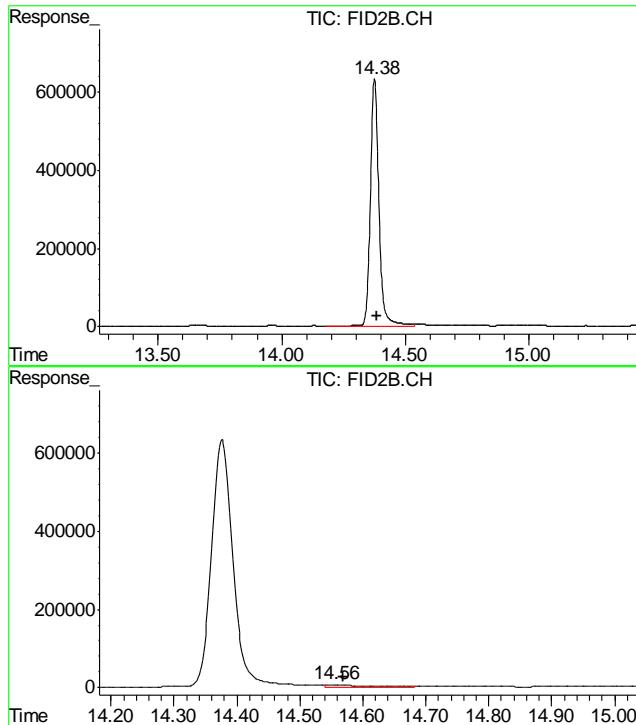
#8 m,p-Xylene

R.T.: 10.481 min
Delta R.T.: -0.011 min
Response: 174040
Conc: 0.10 ug/L



#9 o-Xylene

R.T.: 0.000 min
Exp R.T.: 10.986 min
Response: 0
Conc: N.D.



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

R.T.: 14.376 min
 Delta R.T.: -0.010 min
 Response: 15266238
 Conc: 93.93 %

#11 Naphthalene

R.T.: 14.555 min
 Delta R.T.: -0.013 min
 Response: 233105
 Conc: 1.18 ug/L

11.1.1

Quantitation Report (QT Reviewed)

Signal #1 : Y:\1\DATA\111612\GB18499.D\FID1A.CH Vial: 3
 Signal #2 : Y:\1\DATA\111612\GB18499.D\FID2B.CH
 Acq On : 16 Nov 2012 5:51 pm Operator: StephK
 Sample : MB Inst : GC/MS Ins
 Misc : GC3242,GGB1010,5.000,,100,5,1 Multiplr: 1.00
 IntFile Signal #1: TVH1.E IntFile Signal #2: FB2.E
 Quant Time: Nov 16 18:15:12 2012 Quant Results File: TB868GB868SOIL.RES

Quant Method : C:\MSDCHEM\1...\TB868GB868SOIL.M (Chemstation Integrator)
 Title : 8015B/8021B TVH/BTEX
 Last Update : Fri Nov 16 18:14:55 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	14.38	2895056	92.393	%
10) S 1,2,4-Trichlorobenzene (P)	14.38	15453853	95.085	%

Target Compounds

1) H TVH-Gasoline	7.23	4428413	<MDL	mg/L
4) T Methyl-t-butyl-ether	0.00	0	N.D.	ug/L d
5) T Benzene	0.00	0	N.D.	ug/L d
6) T Toluene	7.68	166985	0.421	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	14.56	187354	0.950	ug/L

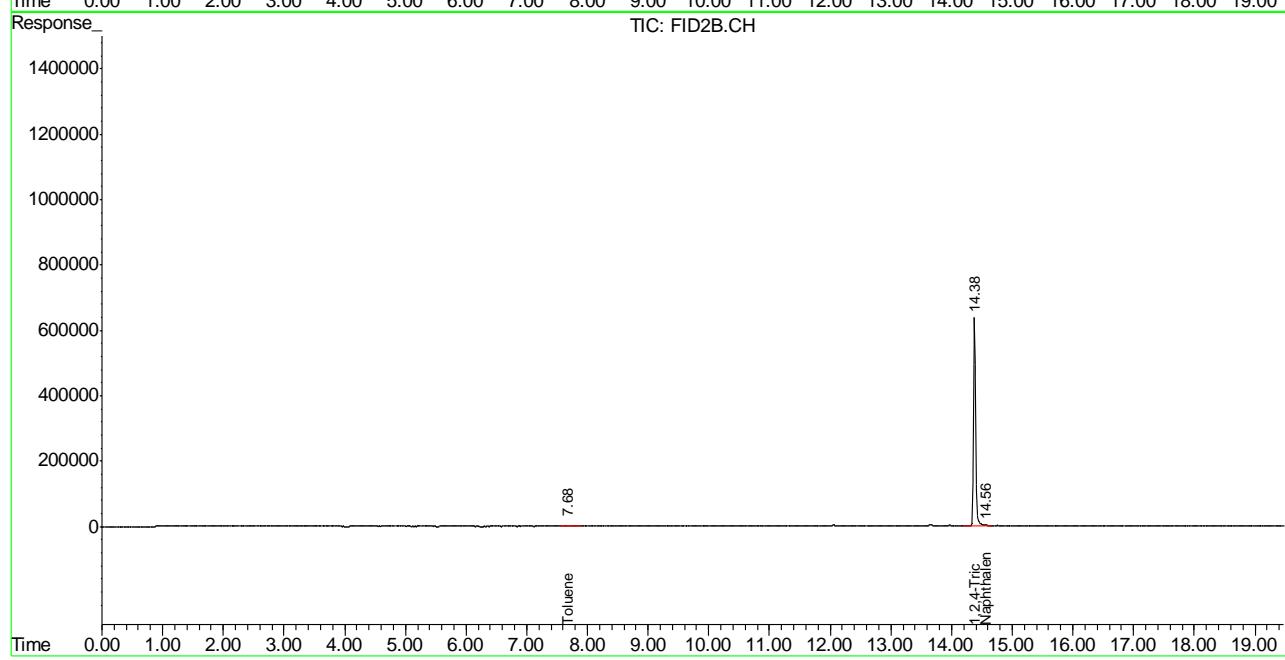
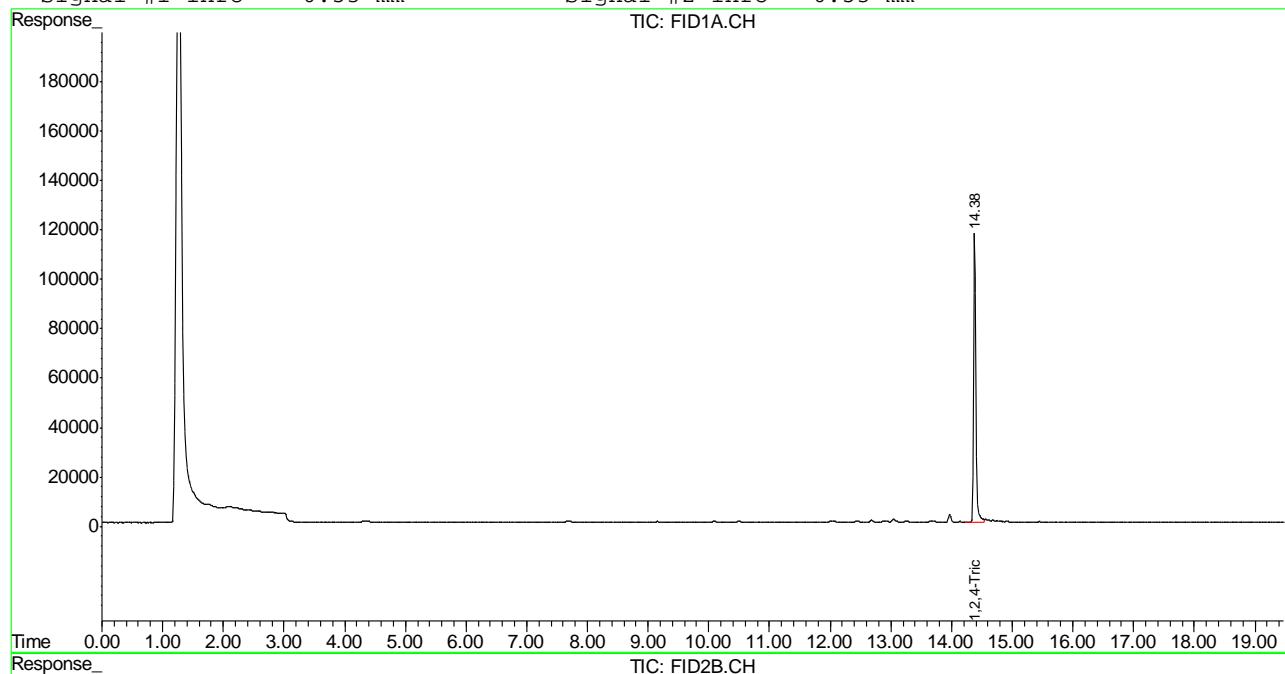
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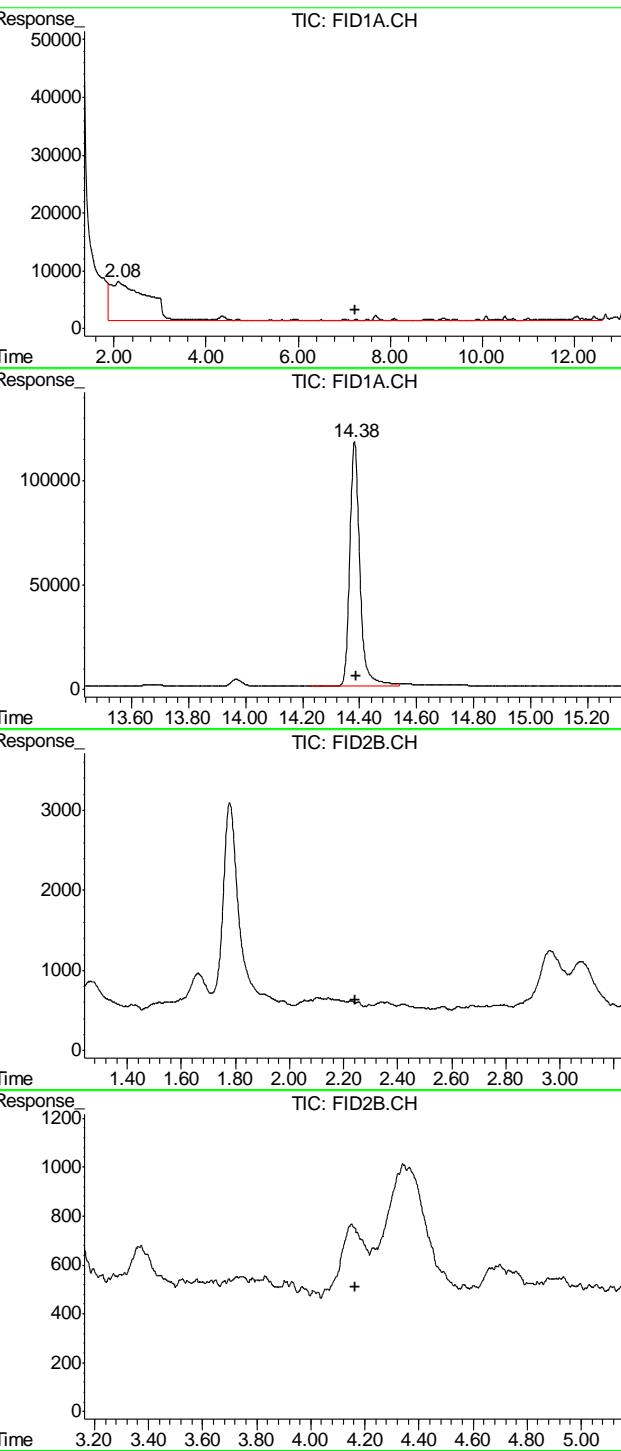
Quantitation Report (QT Reviewed)

Signal #1 : Y:\1\DATA\111612\GB18499.D\FID1A.CH Vial: 3
 Signal #2 : Y:\1\DATA\111612\GB18499.D\FID2B.CH
 Acq On : 16 Nov 2012 5:51 pm Operator: StephK
 Sample : MB Inst : GC/MS Ins
 Misc : GC3242,GGB1010,5.000,,100,5,1 Multiplr: 1.00
 IntFile Signal #1: TVH1.E IntFile Signal #2: FB2.E
 Quant Time: Nov 16 18:14 2012 Quant Results File: TB868GB868SOIL.RES

Quant Method : C:\MSDCHEM\1...\TB868GB868SOIL.M (Chemstation Integrator)
 Title : 8015B/8021B TVH/BTEX
 Last Update : Fri Nov 16 18:14:55 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



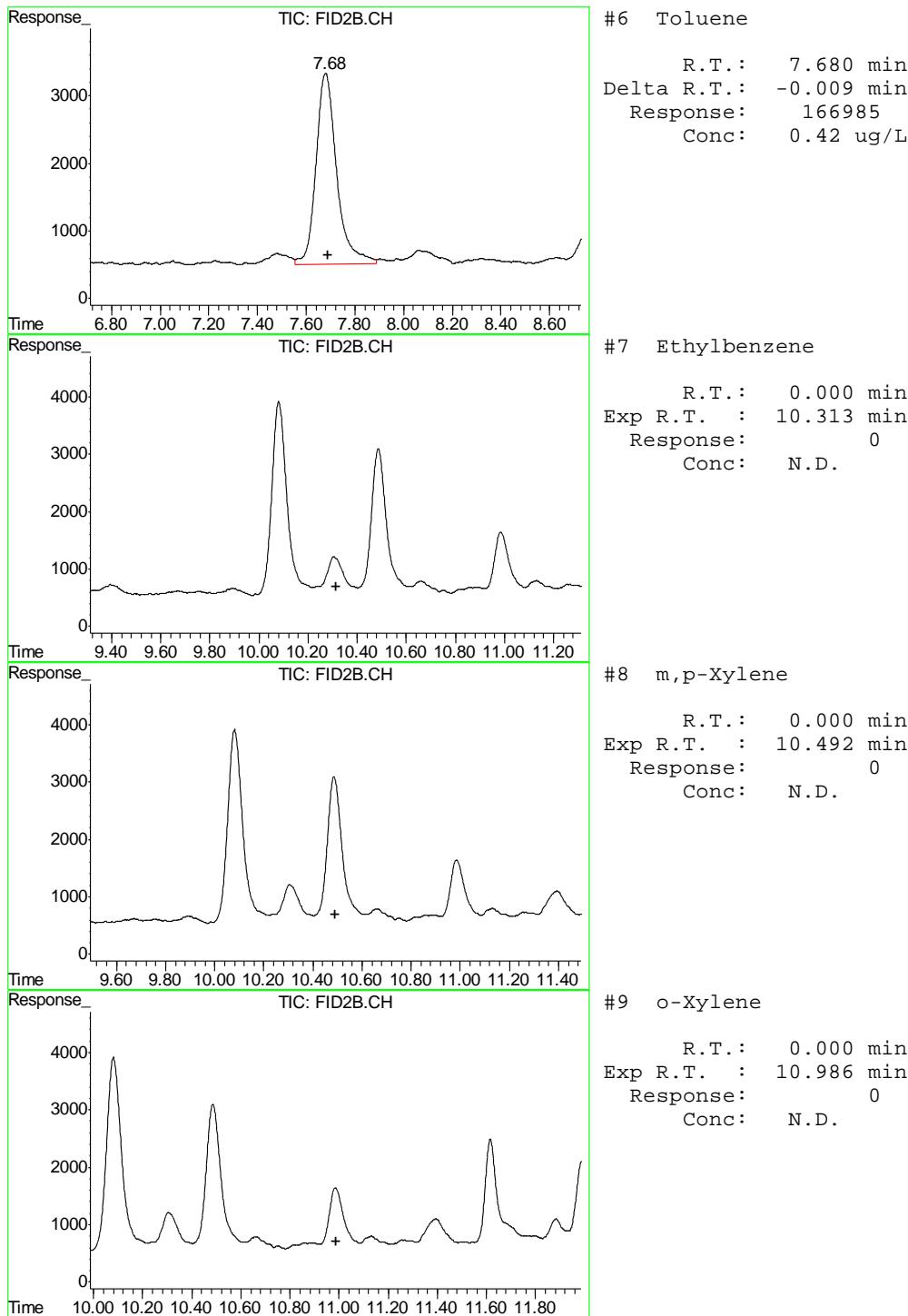


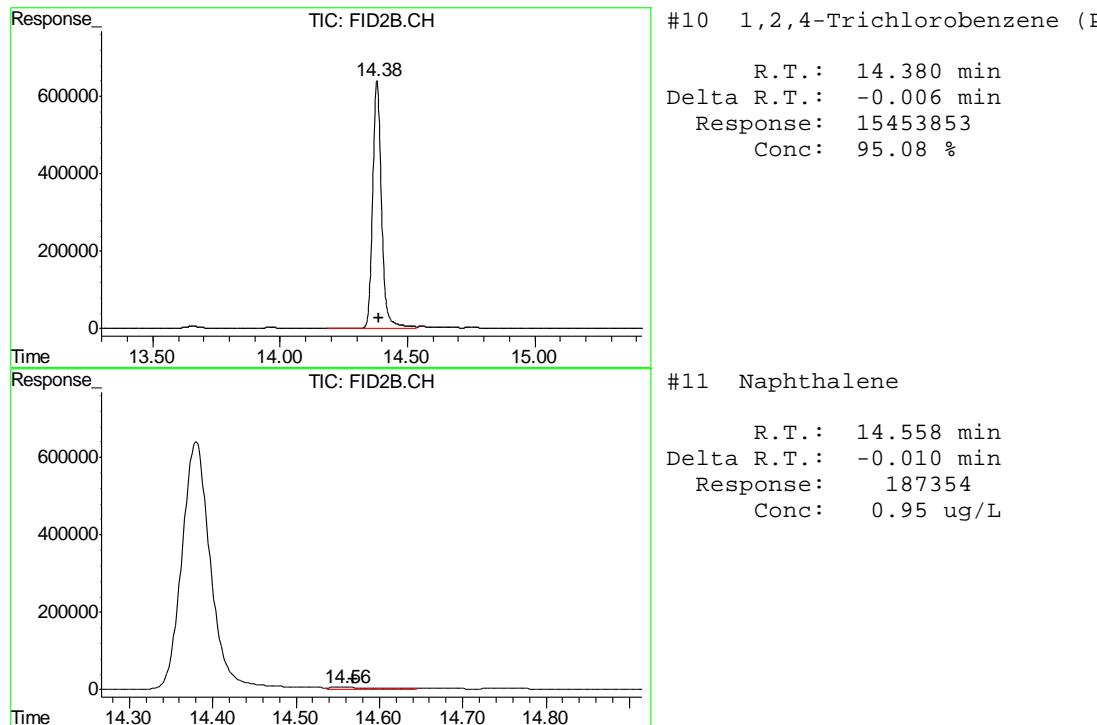
#1 TVH-Gasoline
 R.T.: 7.230 min
 Delta R.T.: 0.000 min
 Response: 4428413
 Conc: N.D.

#2 1,2,4-Trichlorobenzene
 R.T.: 14.382 min
 Delta R.T.: -0.006 min
 Response: 2895056
 Conc: 92.39 %

#4 Methyl-t-butyl-ether
 R.T.: 0.000 min
 Exp R.T. : 2.242 min
 Response: 0
 Conc: N.D.

#5 Benzene
 R.T.: 0.000 min
 Exp R.T. : 4.163 min
 Response: 0
 Conc: N.D.





11.2.1

11



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: D41014

Account: XTOKWR XTO Energy

Project: PCU 296-5A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
OP6979-MB	FD19679.D	1	11/19/12	AV	11/19/12	OP6979	GFD990

The QC reported here applies to the following samples:

Method: SW846-8015B

D41014-1

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

CAS No.	Surrogate Recoveries	Limits
84-15-1	o-Terphenyl	88% 35-130%

Blank Spike Summary

Page 1 of 1

Job Number: D41014

Account: XTOKWR XTO Energy

Project: PCU 296-5A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
OP6979-BS	FD19681.D	1	11/19/12	AV	11/19/12	OP6979	GFD990

The QC reported here applies to the following samples:

Method: SW846-8015B

D41014-1

CAS No.	Compound	Spike mg/kg	BSP mg/kg	BSP %	Limits
	TPH-DRO (C10-C28)	667	689	103	48-130

CAS No.	Surrogate Recoveries	BSP	Limits
84-15-1	o-Terphenyl	95%	35-130%

* = Outside of Control Limits.

12.2.1
12

Matrix Spike/Matrix Spike Duplicate Summary

Page 1 of 1

Job Number: D41014

Account: XTOKWR XTO Energy

Project: PCU 296-5A

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
OP6979-MS	FD19683.D	1	11/19/12	AV	11/19/12	OP6979	GFD990
OP6979-MSD	FD19685.D	1	11/19/12	AV	11/19/12	OP6979	GFD990
D41044-9	FD19687.D	1	11/19/12	AV	11/19/12	OP6979	GFD990

The QC reported here applies to the following samples:

Method: SW846-8015B

D41014-1

CAS No.	Compound	D41044-9		Spike	MS	MS	MSD	MSD	RPD	Limits Rec/RPD
		mg/kg	Q	mg/kg	mg/kg	%	mg/kg	%		
	TPH-DRO (C10-C28)	80.3		738	764	93	734	89	4	20-168/30
CAS No.	Surrogate Recoveries	MS		MSD		D41044-9	Limits			
84-15-1	o-Terphenyl	80%		80%		65%	35-130%			

* = Outside of Control Limits.

12.3.1
12



GC Semi-volatiles

Raw Data

Quantitation Report (QT Reviewed)

Data File : C:\MSDCHEM\2\DATA\2012\NOV\FD111912\FD19691.D Vial: 9
 Acq On : 11-19-2012 05:37:12 PM Operator: ashleyv
 Sample : D41014-1 Inst : FID5
 Misc : OP6979,GFD990,30.01,,,2,1 Multiplr: 1.00
 IntFile : autoint1.e
 Quant Time: Nov 20 15:06:13 2012 Quant Results File: DRO-GFD982F.RES

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

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

Compound	R.T.	Response	Conc	Units
<hr/>				
System Monitoring Compounds				
1) S O-Terphenyl	8.95	43488217	785.974	mg/L
<hr/>				
Target Compounds				
2) H TPH-DRO (c10-c28)	6.89	12147235	319.869	mg/L

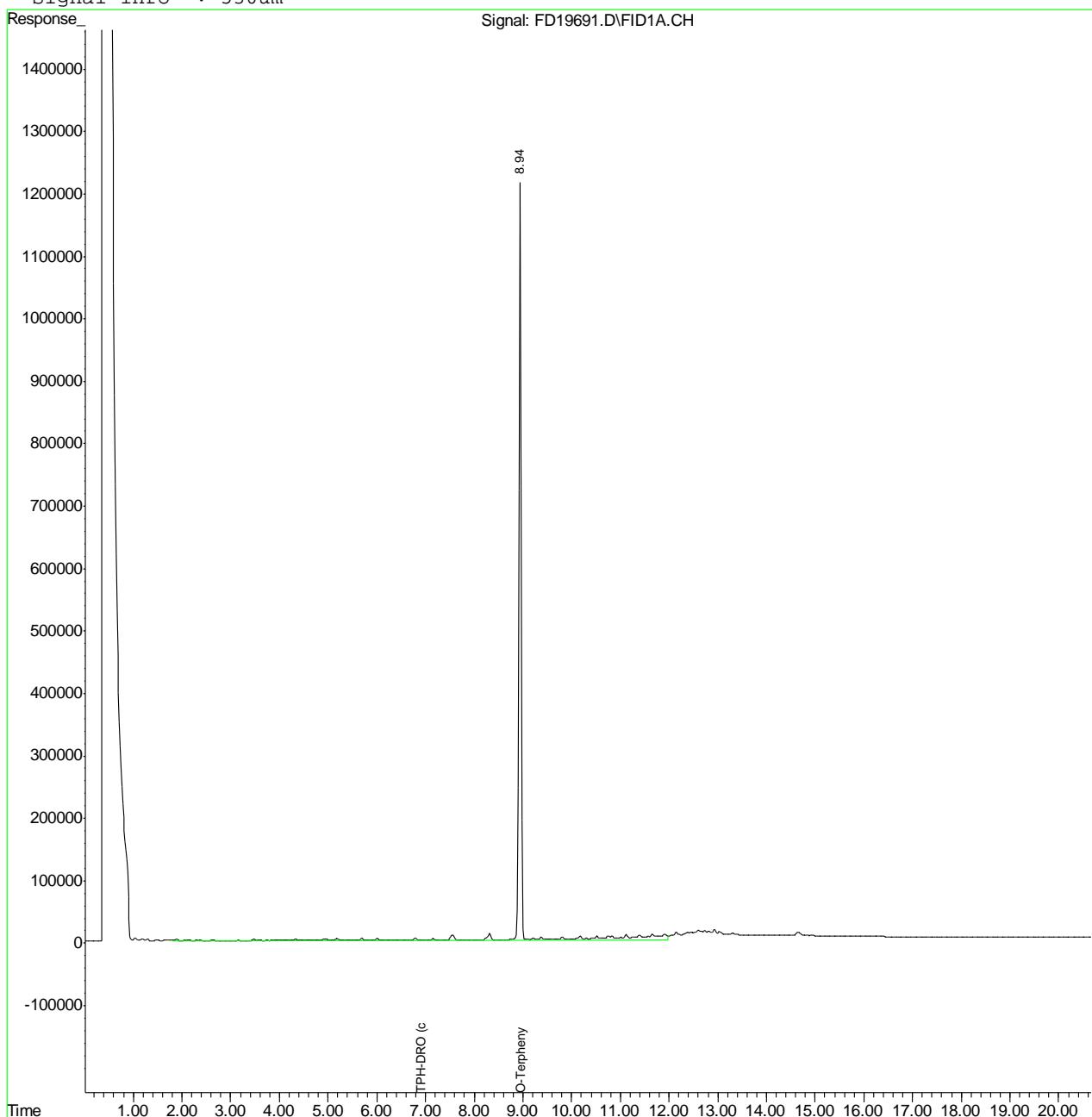
(f)=RT Delta > 1/2 Window (m)=manual int.
 FD19691.D DRO-GFD982F.M Wed Nov 21 09:05:45 2012 GC

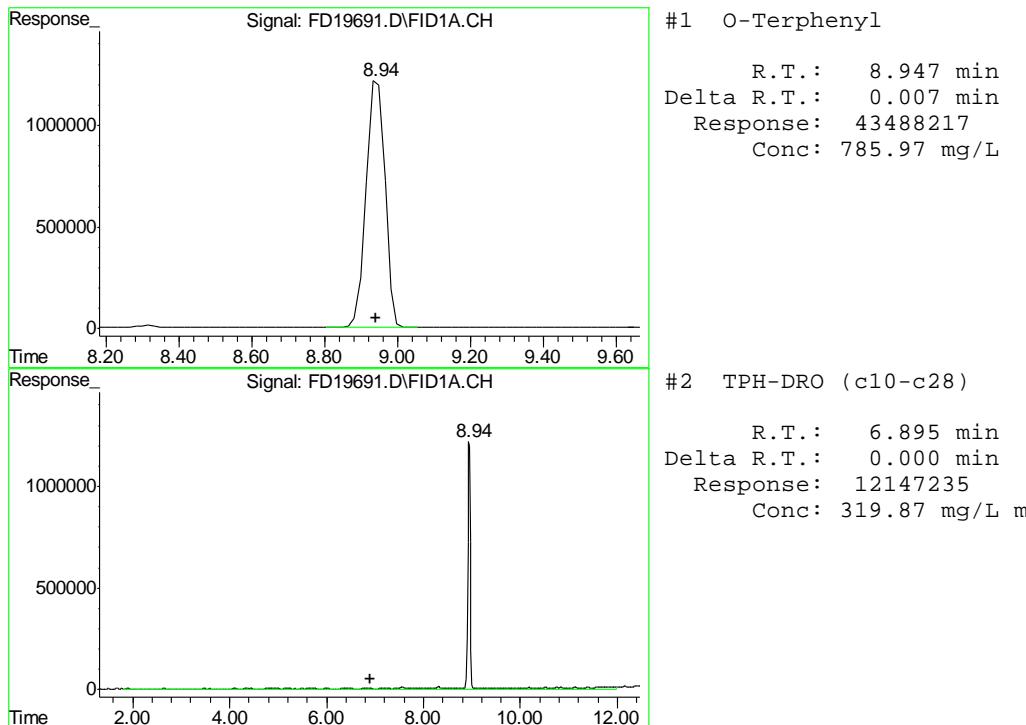
Quantitation Report (QT Reviewed)

Data File : C:\MSDCHEM\2\DATA\2012\NOV\FD111912\FD19691.D Vial: 9
 Acq On : 11-19-2012 05:37:12 PM Operator: ashleyv
 Sample : D41014-1 Inst : FID5
 Misc : OP6979,GFD990,30.01,,,2,1 Multiplr: 1.00
 IntFile : autoint1.e
 Quant Time: Nov 20 15:06 2012 Quant Results File: DRO-GFD982F.RES

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

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





Quantitation Report (QT Reviewed)

Data File : C:\MSDCHEM\2\DATA\2012\NOV\FD111912\FD19679.D Vial: 3
 Acq On : 11-19-2012 02:55:37 PM Operator: ashleyv
 Sample : OP6979-MB Inst : FID5
 Misc : OP6979,GFD990,30.00,,,2,1 Multiplr: 1.00
 IntFile : autoint1.e
 Quant Time: Nov 20 15:02:12 2012 Quant Results File: DRO-GFD982F.RES

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

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

Compound	R.T.	Response	Conc Units
<hr/>			
System Monitoring Compounds			
1) S O-Terphenyl	8.97	48861805	883.092 mg/L
<hr/>			
Target Compounds			
2) H TPH-DRO (c10-c28)	6.89	2183565	57.499 mg/L

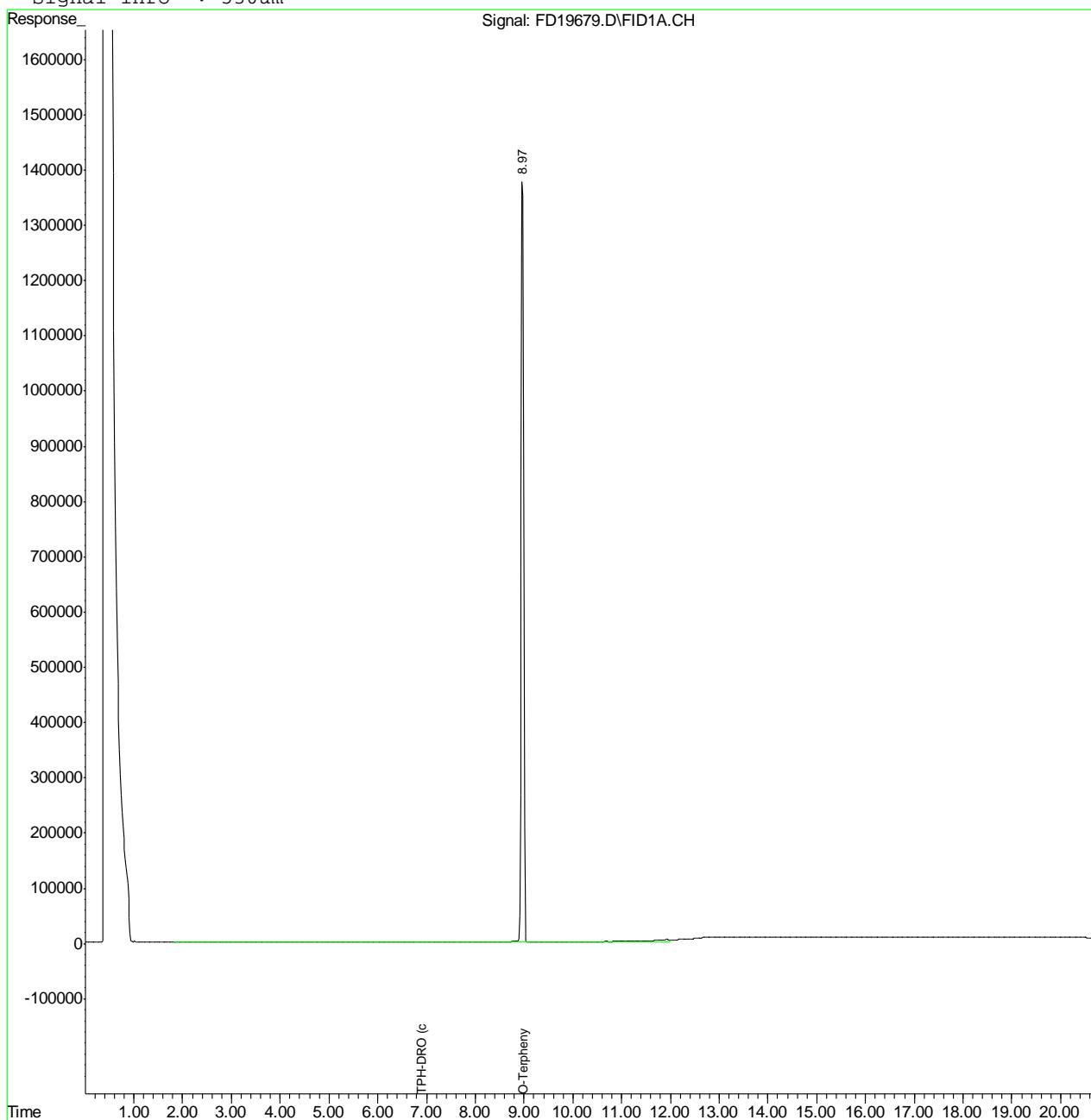
(f)=RT Delta > 1/2 Window (m)=manual int.
 FD19679.D DRO-GFD982F.M Wed Nov 21 09:05:39 2012 GC

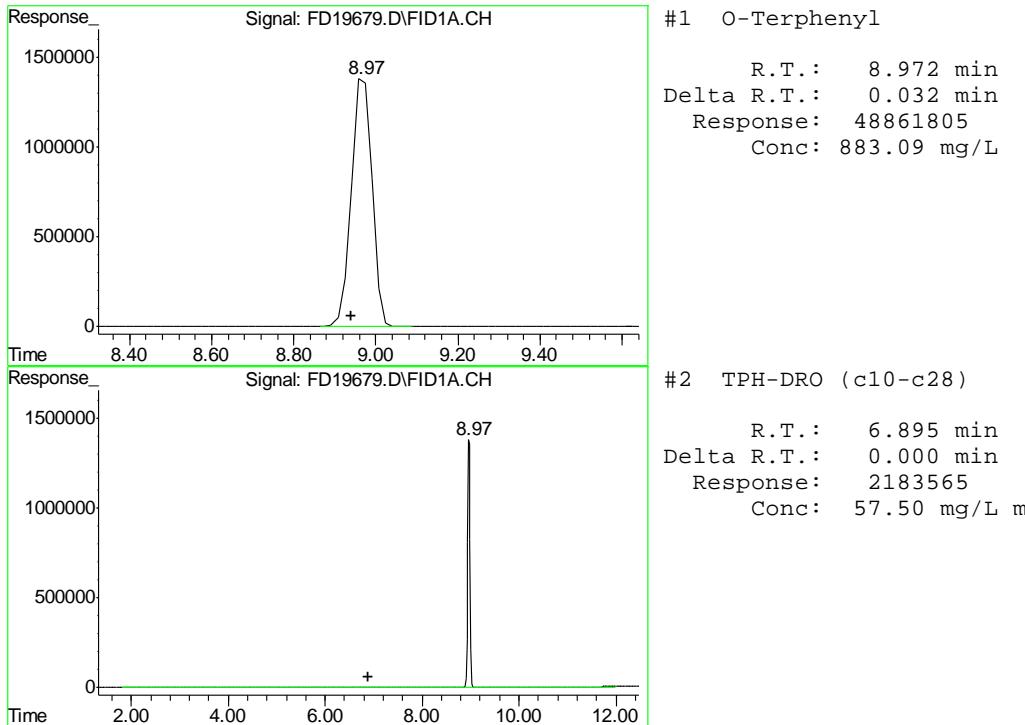
Quantitation Report (QT Reviewed)

Data File : C:\MSDCHEM\2\DATA\2012\NOV\FD111912\FD19679.D Vial: 3
 Acq On : 11-19-2012 02:55:37 PM Operator: ashleyv
 Sample : OP6979-MB Inst : FID5
 Misc : OP6979,GFD990,30.00,,,2,1 Multiplr: 1.00
 IntFile : autoint1.e
 Quant Time: Nov 20 15:02 2012 Quant Results File: DRO-GFD982F.RES

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

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





13.2.1

13



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: D41014
Account: XTOKRWR - XTO Energy
Project: PCU 296-5A

QC Batch ID: MP8913
Matrix Type: SOLID

Methods: SW846 6010C
Units: mg/kg

Prep Date:

11/19/12

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	0.020	<1.0
Beryllium	1.0	.13	.15		
Boron	5.0	.1	.06		
Cadmium	1.0	.06	.036	-0.020	<1.0
Calcium	40	.54	9		
Chromium	1.0	.03	.03	0.010	<1.0
Cobalt	0.50	.04	.07		
Copper	1.0	.12	.15	-0.070	<1.0
Iron	7.0	.12	.87		
Lead	5.0	.19	.24	-0.33	<5.0
Lithium	0.20	.05	.054		
Magnesium	20	.65	.98		
Manganese	0.50	.12	.022		
Molybdenum	1.0	.21	.08		
Nickel	3.0	.05	.026	0.0	<3.0
Phosphorus	10	1.4	1.9		
Potassium	200	6.1	7		
Selenium	5.0	.48	.36	-0.21	<5.0
Silicon	5.0	.29	.37		
Silver	3.0	.04	.06	-0.040	<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	-0.18	<3.0

Associated samples MP8913: D41014-1

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

BLANK RESULTS SUMMARY
Part 2 - Method Blanks

Login Number: D41014
Account: XTOKRWR - XTO Energy
Project: PCU 296-5A

QC Batch ID: MP8913
Matrix Type: SOLID

Methods: SW846 6010C
Units: mg/kg

Prep Date:

Metal

(anr) Analyte not requested

MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D41014
 Account: XTOKRWR - XTO Energy
 Project: PCU 296-5A

QC Batch ID: MP8913
 Matrix Type: SOLID

Methods: SW846 6010C
 Units: mg/kg

Prep Date:

11/19/12

Metal	D41013-1 Original MS	Spikelot ICPALL2	% Rec	QC Limits
Aluminum	anr			
Antimony	anr			
Arsenic	anr			
Barium	1000	1310	244	127.2(a) 75-125
Beryllium	anr			
Boron				
Cadmium	0.0	49.1	60.9	80.6 75-125
Calcium				
Chromium	34.1	80.8	60.9	76.7 75-125
Cobalt	anr			
Copper	12.6	63.5	60.9	83.5 75-125
Iron	anr			
Lead	11.5	112	122	82.5 75-125
Lithium				
Magnesium				
Manganese	anr			
Molybdenum				
Nickel	17.5	61.9	60.9	72.9N(b) 75-125
Phosphorus	anr			
Potassium				
Selenium	0.77	99.9	122	81.4 75-125
Silicon				
Silver	0.18	18.1	24.4	73.5N(b) 75-125
Sodium				
Strontium				
Thallium	anr			
Tin				
Titanium				
Uranium				
Vanadium	anr			
Zinc	46.7	96.6	60.9	81.9 75-125

Associated samples MP8913: D41014-1

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

MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D41014
Account: XTOKRWR - XTO Energy
Project: PCU 296-5A

QC Batch ID: MP8913
Matrix Type: SOLID

Methods: SW846 6010C
Units: mg/kg

Prep Date:

Metal

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

14.1.2
14

MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D41014
 Account: XTOKRWR - XTO Energy
 Project: PCU 296-5A

QC Batch ID: MP8913
 Matrix Type: SOLID

Methods: SW846 6010C
 Units: mg/kg

Prep Date:

11/19/12

Metal	D41013-1 Original	MSD	Spikelot ICPALL2	% Rec	MSD RPD	QC Limit
Aluminum	anr					
Antimony	anr					
Arsenic	anr					
Barium	1000	1220	251	87.5	7.1	20
Beryllium	anr					
Boron						
Cadmium	0.0	50.0	62.8	79.6	1.8	20
Calcium						
Chromium	34.1	77.5	62.8	69.1N(a)	4.2	20
Cobalt	anr					
Copper	12.6	64.7	62.8	82.9	1.9	20
Iron	anr					
Lead	11.5	113	126	80.8	0.9	20
Lithium						
Magnesium						
Manganese	anr					
Molybdenum						
Nickel	17.5	62.3	62.8	71.3N(a)	0.6	20
Phosphorus	anr					
Potassium						
Selenium	0.77	101	126	79.8	1.1	20
Silicon						
Silver	0.18	18.5	25.1	72.9N(a)	2.2	20
Sodium						
Strontium						
Thallium	anr					
Tin						
Titanium						
Uranium						
Vanadium	anr					
Zinc	46.7	96.1	62.8	78.6	0.5	20

Associated samples MP8913: D41014-1

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

MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D41014
Account: XTOKRWR - XTO Energy
Project: PCU 296-5A

QC Batch ID: MP8913
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: D41014
 Account: XTOKRWR - XTO Energy
 Project: PCU 296-5A

QC Batch ID: MP8913
 Matrix Type: SOLID

Methods: SW846 6010C
 Units: mg/kg

Prep Date:

11/19/12

Metal	BSP Result	Spikelot ICPALL2	% Rec	QC Limits
Aluminum	anr			
Antimony	anr			
Arsenic	anr			
Barium	190	200	95.0	80-120
Beryllium	anr			
Boron				
Cadmium	45.1	50	90.2	80-120
Calcium				
Chromium	47.7	50	95.4	80-120
Cobalt	anr			
Copper	43.0	50	86.0	80-120
Iron	anr			
Lead	96.2	100	96.2	80-120
Lithium				
Magnesium				
Manganese	anr			
Molybdenum				
Nickel	45.7	50	91.4	80-120
Phosphorus	anr			
Potassium				
Selenium	91.1	100	91.1	80-120
Silicon				
Silver	16.8	20	84.0	80-120
Sodium				
Strontium				
Thallium	anr			
Tin				
Titanium				
Uranium				
Vanadium	anr			
Zinc	48.3	50	96.6	80-120

Associated samples MP8913: D41014-1

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

SPIKE BLANK AND LAB CONTROL SAMPLE SUMMARY

Login Number: D41014
Account: XTOKRWR - XTO Energy
Project: PCU 296-5A

QC Batch ID: MP8913
Matrix Type: SOLID

Methods: SW846 6010C
Units: mg/kg

Prep Date:

Metal

(anr) Analyte not requested

SERIAL DILUTION RESULTS SUMMARY

Login Number: D41014
 Account: XTOKRWR - XTO Energy
 Project: PCU 296-5A

QC Batch ID: MP8913
 Matrix Type: SOLID

Methods: SW846 6010C
 Units: ug/l

Prep Date:

11/19/12

Metal	D41013-1 Original	SDL 1:5	%DIF	QC Limits
Aluminum	anr			
Antimony	anr			
Arsenic	anr			
Barium	8220	9560	16.3*(a)	0-10
Beryllium	anr			
Boron				
Cadmium	0.00	0.00	NC	0-10
Calcium				
Chromium	280	326	16.4*(a)	0-10
Cobalt	anr			
Copper	103	106	2.3	0-10
Iron	anr			
Lead	94.0	92.0	2.1	0-10
Lithium				
Magnesium				
Manganese	anr			
Molybdenum				
Nickel	144	173	19.8*(a)	0-10
Phosphorus	anr			
Potassium				
Selenium	6.30	0.00	100.0(b)	0-10
Silicon				
Silver	1.50	0.00	100.0(b)	0-10
Sodium				
Strontium				
Thallium	anr			
Tin				
Titanium				
Uranium				
Vanadium	anr			
Zinc	383	458	19.5*(a)	0-10

Associated samples MP8913: D41014-1

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

SERIAL DILUTION RESULTS SUMMARY

Login Number: D41014
Account: XTOKRWR - XTO Energy
Project: PCU 296-5A

QC Batch ID: MP8913
Matrix Type: SOLID

Methods: SW846 6010C
Units: ug/l

Prep Date:

Metal

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

BLANK RESULTS SUMMARY
Part 2 - Method Blanks

Login Number: D41014
Account: XTOKRWR - XTO Energy
Project: PCU 296-5A

QC Batch ID: MP8914
Matrix Type: SOLID

Methods: SW846 6020A
Units: mg/kg

Prep Date:

11/19/12

Metal	RL	IDL	MDL	MB raw	final
Aluminum	25	.22	.31		
Antimony	0.20	.0018	.0075		
Arsenic	0.10	.006	.06	0.0076	<0.10
Barium	1.0	.0065	.037		
Beryllium	0.10	.016	.09		
Boron	20	1.2	1.2		
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		
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 MP8914: D41014-1

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: D41014
 Account: XTOKRWR - XTO Energy
 Project: PCU 296-5A

QC Batch ID: MP8914
 Matrix Type: SOLID

Methods: SW846 6020A
 Units: mg/kg

Prep Date:

11/19/12

Metal	D41013-1 Original MS	Spikelot ICPALL2	% Rec	QC Limits
Aluminum				
Antimony				
Arsenic	8.1	128	122	98.4 75-125
Barium				
Beryllium				
Boron				
Calcium				
Cobalt				
Iron				
Lead	anr			
Magnesium				
Manganese				
Molybdenum	anr			
Potassium	anr			
Selenium	anr			
Silver				
Sodium	anr			
Strontium				
Thallium				
Tin				
Titanium				
Uranium				
Vanadium				

Associated samples MP8914: D41014-1

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

14.2.2
14

MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D41014
 Account: XTOKRWR - XTO Energy
 Project: PCU 296-5A

QC Batch ID: MP8914
 Matrix Type: SOLID

Methods: SW846 6020A
 Units: mg/kg

Prep Date:

11/19/12

Metal	D41013-1 Original	MSD	Spikelot ICPALL2	% Rec	MSD RPD	QC Limit
Aluminum						
Antimony						
Arsenic	8.1	130	126	97.0	1.6	20
Barium						
Beryllium						
Boron						
Calcium						
Cobalt						
Iron						
Lead	anr					
Magnesium						
Manganese						
Molybdenum	anr					
Potassium	anr					
Selenium	anr					
Silver						
Sodium	anr					
Strontium						
Thallium						
Tin						
Titanium						
Uranium						
Vanadium						

Associated samples MP8914: D41014-1

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

14.2.2
14

SPIKE BLANK AND LAB CONTROL SAMPLE SUMMARY

Login Number: D41014
 Account: XTOKRWR - XTO Energy
 Project: PCU 296-5A

QC Batch ID: MP8914
 Matrix Type: SOLID

Methods: SW846 6020A
 Units: mg/kg

Prep Date:

11/19/12

Metal	BSP Result	Spikelot ICPALL2	% Rec	QC Limits
Aluminum				
Antimony				
Arsenic	100	100	100.0	80-120
Barium				
Beryllium				
Boron				
Calcium				
Chromium	anr			
Cobalt				
Copper	anr			
Iron				
Lead	anr			
Magnesium				
Manganese				
Molybdenum	anr			
Nickel	anr			
Potassium	anr			
Selenium	anr			
Silver				
Sodium	anr			
Strontium				
Thallium				
Tin				
Titanium				
Uranium				
Vanadium				
Zinc	anr			

Associated samples MP8914: D41014-1

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: D41014
 Account: XTOKRWR - XTO Energy
 Project: PCU 296-5A

QC Batch ID: MP8914
 Matrix Type: SOLID

Methods: SW846 6020A
 Units: ug/l

Prep Date:

11/19/12

Metal	D41013-1 Original	SDL 5:25	%DIF	QC Limits
Aluminum				
Antimony				
Arsenic	66.8	73.5	9.9	0-10
Barium				
Beryllium				
Boron				
Calcium				
Cobalt				
Iron				
Lead		anr		
Magnesium				
Manganese				
Molybdenum		anr		
Potassium		anr		
Selenium		anr		
Silver				
Sodium		anr		
Strontium				
Thallium				
Tin				
Titanium				
Uranium				
Vanadium				

Associated samples MP8914: D41014-1

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: D41014
Account: XTOKRWR - XTO Energy
Project: PCU 296-5A

QC Batch ID: MP8915
Matrix Type: AQUEOUS

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

Prep Date:

11/19/12

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	22.5	<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	38.0	<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	-290	<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 MP8915: D41014-1A

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

BLANK RESULTS SUMMARY
Part 2 - Method Blanks

Login Number: D41014
Account: XTOKRWR - XTO Energy
Project: PCU 296-5A

QC Batch ID: MP8915
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: D41014
 Account: XTOKRWR - XTO Energy
 Project: PCU 296-5A

QC Batch ID: MP8915
 Matrix Type: AQUEOUS

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

Prep Date: 11/19/12

Metal	D41042-1A Original MS	Spikelot ICPALL2	% Rec	QC Limits
Aluminum				
Antimony				
Arsenic				
Barium				
Beryllium				
Boron				
Cadmium				
Calcium	194000	332000	125000	110.4
Chromium				
Cobalt				
Copper				
Iron				
Lead				
Lithium				
Magnesium	99900	226000	125000	100.9
Manganese				
Molybdenum				
Nickel				
Phosphorus				
Potassium				
Selenium				
Silicon				
Silver				
Sodium	660000	765000	125000	84.0
Strontium				
Thallium				
Tin				
Titanium				
Uranium				
Vanadium				
Zinc				

Associated samples MP8915: D41014-1A

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

MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D41014
Account: XTOKRWR - XTO Energy
Project: PCU 296-5A

QC Batch ID: MP8915
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: D41014
 Account: XTOKRWR - XTO Energy
 Project: PCU 296-5A

QC Batch ID: MP8915
 Matrix Type: AQUEOUS

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

Prep Date:

11/19/12

Metal	D41042-1A Original MSD	Spikelot ICPALL2	MSD % Rec	MSD RPD	QC Limit
Aluminum					
Antimony					
Arsenic					
Barium					
Beryllium					
Boron					
Cadmium					
Calcium	194000	337000	125000	114.4	1.5
Chromium					
Cobalt					
Copper					
Iron					
Lead					
Lithium					
Magnesium	99900	221000	125000	96.9	2.2
Manganese					
Molybdenum					
Nickel					
Phosphorus					
Potassium					
Selenium					
Silicon					
Silver					
Sodium	660000	746000	125000	68.8 (a)	2.5
Strontium					
Thallium					
Tin					
Titanium					
Uranium					
Vanadium					
Zinc					

Associated samples MP8915: D41014-1A

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

MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D41014
Account: XTOKRWR - XTO Energy
Project: PCU 296-5A

QC Batch ID: MP8915
Matrix Type: AQUEOUS

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

Prep Date:

Metal

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

14.3.2
14

SPIKE BLANK AND LAB CONTROL SAMPLE SUMMARY

Login Number: D41014
 Account: XTOKWR - XTO Energy
 Project: PCU 296-5A

QC Batch ID: MP8915
 Matrix Type: AQUEOUS

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

Prep Date: 11/19/12

Metal	BSP Result	Spikelot ICPALL2	% Rec	QC Limits
Aluminum				
Antimony				
Arsenic				
Barium				
Beryllium				
Boron				
Cadmium				
Calcium	141000	125000	112.8	80-120
Chromium				
Cobalt				
Copper				
Iron				
Lead				
Lithium				
Magnesium	126000	125000	100.8	80-120
Manganese				
Molybdenum				
Nickel				
Phosphorus				
Potassium				
Selenium				
Silicon				
Silver				
Sodium	122000	125000	97.6	80-120
Strontium				
Thallium				
Tin				
Titanium				
Uranium				
Vanadium				
Zinc				

Associated samples MP8915: D41014-1A

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

SPIKE BLANK AND LAB CONTROL SAMPLE SUMMARY

Login Number: D41014
Account: XTOKRWR - XTO Energy
Project: PCU 296-5A

QC Batch ID: MP8915
Matrix Type: AQUEOUS

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

Prep Date:

Metal

(anr) Analyte not requested

SERIAL DILUTION RESULTS SUMMARY

Login Number: D41014
 Account: XTOKRWR - XTO Energy
 Project: PCU 296-5A

QC Batch ID: MP8915
 Matrix Type: AQUEOUS

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

Prep Date: 11/19/12

Metal	D41042-1A	Original	SDL 1:5	%DIF	QC Limits
Aluminum					
Antimony					
Arsenic					
Barium					
Beryllium					
Boron					
Cadmium					
Calcium	38800	40600	4.9		0-10
Chromium					
Cobalt					
Copper					
Iron					
Lead					
Lithium					
Magnesium	20000	22200	11.3*(a)		0-10
Manganese					
Molybdenum					
Nickel					
Phosphorus					
Potassium					
Selenium					
Silicon					
Silver					
Sodium	132000	147000	11.4*(a)		0-10
Strontium					
Thallium					
Tin					
Titanium					
Uranium					
Vanadium					
Zinc					

Associated samples MP8915: D41014-1A

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

SERIAL DILUTION RESULTS SUMMARY

Login Number: D41014
Account: XTOKRWR - XTO Energy
Project: PCU 296-5A

QC Batch ID: MP8915
Matrix Type: AQUEOUS

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

Prep Date:

Metal

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

BLANK RESULTS SUMMARY
Part 2 - Method Blanks

Login Number: D41014
Account: XTOKRWR - XTO Energy
Project: PCU 296-5A

QC Batch ID: MP8936
Matrix Type: SOLID

Methods: SW846 7471B
Units: mg/kg

Prep Date:

11/27/12

Metal	RL	IDL	MDL	MB raw	final
Mercury	0.083	.00088	.00075	0.0013	<0.083

Associated samples MP8936: D41014-1

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

14.4.1
14

MATRIX SPIKE AND DUPLICATE RESULTS SUMMARY

Login Number: D41014
Account: XTOKWR - XTO Energy
Project: PCU 296-5A

QC Batch ID: MP8936
Matrix Type: SOLID

Methods: SW846 7471B
Units: mg/kg

Prep Date:

11/27/12

Metal	D40988-1 Original MS	Spikelot HGWSR1	QC % Rec	QC Limits
Mercury	0.029	0.39	0.369	97.8 75-125

Associated samples MP8936: D41014-1

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: D41014
Account: XTOKRWR - XTO Energy
Project: PCU 296-5A

QC Batch ID: MP8936
Matrix Type: SOLID

Methods: SW846 7471B
Units: mg/kg

Prep Date:

11/27/12

Metal	D40988-1 Original	MSD	Spikelot HGWSR1	MSD % Rec	RPD	QC Limit
Mercury	0.029	0.39	0.382	94.6	0.0	20

Associated samples MP8936: D41014-1

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: D41014
Account: XTOKRWR - XTO Energy
Project: PCU 296-5A

QC Batch ID: MP8936
Matrix Type: SOLID

Methods: SW846 7471B
Units: mg/kg

Prep Date: 11/27/12

Metal	BSP Result	Spikelot HGWSR1	QC % Rec	QC Limits
Mercury	0.33	0.333	99.0	80-120

Associated samples MP8936: D41014-1

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

14.4.3
14



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: D41014
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	GP8709/GN17735	1.0	0.0	mg/kg	176.0	166	94.0	80-120%
Specific Conductivity	GP8712/GN17742			umhos/cm	9991	9990	100.0	90-110%
pH	GN17719			su	8.00	8.00	100.0	99.3-100.7%

Associated Samples:

Batch GP8709: D41014-1

Batch GP8712: D41014-1

Batch GN17719: D41014-1

(*) Outside of QC limits

DUPLICATE RESULTS SUMMARY
GENERAL CHEMISTRY

Login Number: D41014
Account: XTOKWR - XTO Energy
Project: PCU 296-5A

Analyte	Batch ID	QC Sample	Units	Original Result	DUP Result	RPD	QC Limits
Chromium, Hexavalent Redox Potential Vs H2	GP8709/GN17735 GN17722	D40985-1 D41014-1	mg/kg mv	0.0 94.2	0.0 93.7	0.0 0.5	0-20% 0-20%

Associated Samples:

Batch GP8709: D41014-1

Batch GN17722: D41014-1

(*) Outside of QC limits

MATRIX SPIKE RESULTS SUMMARY
GENERAL CHEMISTRY

Login Number: D41014
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	GP8709/GN17735	D40985-1	mg/kg	0.0	40.0	39.8	99.5	75-125%

Associated Samples:

Batch GP8709: D41014-1

(*) Outside of QC limits

(N) Matrix Spike Rec. outside of QC limits

MATRIX SPIKE DUPLICATE RESULTS SUMMARY
GENERAL CHEMISTRY

Login Number: D41014
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	GP8709/GN17735	D40985-1	mg/kg	0.0	40.0	40.9	2.7	20%

Associated Samples:

Batch GP8709: D41014-1

(*) Outside of QC limits

(N) Matrix Spike Rec. outside of QC limits