908 North Temperance Ave. ∇ Clovis, CA 93611 ∇ Phone 559-275-2175 ∇ Fax 559-275-4422

Certification Number: CA1312
NELAP Certification number: 05233CA
DoD-ELAP Certificate number: ADE-1410

Data Validatable Report

December 12, 2011

Environet, Inc.
650 Iwilei Road, Suite 204
Honolulu, Hawaii 96817

Attn: Stacey Fineran

Title: Report of Data: Case 66102

Project: 1022-024 LTM Red Hill Bulk Fuel Storage Facility

Contract #: Prime contract # for DoD: N62742-08-D-1930, CTO HC21

Dear Ms. Fineran:

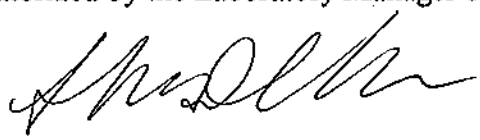
Samples were received October 26, 2011, in good condition. Written results for the requested analyses are provided on this December 12, 2011.

Results in this report apply to the samples analyzed in accordance with the chain of custody document. This analytical report must be reproduced in its entirety.

The MADEP-EPH and VPH analyses were subcontracted to Gulf coast Analytical Laboratories, Inc.

If you have any questions or require further information, please contact your APPL Project Manager, Diane Anderson, at your convenience. Thank you for choosing APPL, Inc.

I certify that this data package is in compliance with the terms and conditions of the contract, both technically and for completeness, for other than the conditions detailed above. These test results meet all requirements of NELAC and DoD QSM. Release of the hard copy has been authorized by the Laboratory Manager or her designee, as verified by the following signature.



Sharon Dehmlow, Laboratory Director
APPL, Inc.

SD/rp
Enclosure
cc: File

Number of pages in this report: ___

908 North Temperance Ave. ∇ Clovis, CA 93611 ∇ Phone 559-275-2175 ∇ Fax 559-275-4422

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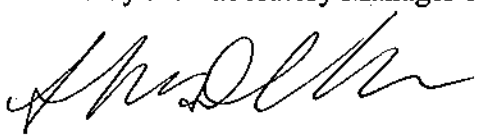


Sharon Dehmlow, Laboratory Director
APPL, Inc.

SD/rp
Enclosure
cc: File

Number of pages in this report: 483

Data Validation Package
for
LTM Red Hill Bulk Fuel Storage Facility
SDG 66102
TABLE OF CONTENTS

LABORATORY NAME: *APPL, Inc.*

SAMPLE RECEIPT INFORMATION

Sample receipt information

ARF: 66102

Project: LTM Red Hill Bulk Fuel Storage Facility

Sample Receipt Information:

The samples were received on October 26, 2011, at 2.5°C, 3.0°C, and 3.0°C. The samples were assigned Analytical Request Form (ARF) number 66102. The sample numbers and requested analyses were compared to the chains of custody and email communications. Four bottles arrived broken for sample ES047; the client was notified. No other exception was encountered.

Sample Table

CLIENT ID	APPL ID	Matrix	Date Sampled	Date Received
ES046	AY49333	WATER	10/24/2011	10/26/2011
ES047	AY49334	WATER	10/24/2011	10/26/2011
ES048	AY49335	WATER	10/24/2011	10/26/2011
ES049	AY49336	WATER	10/24/2011	10/26/2011

Samples and blanks were screened for J-value responses between the detection limit (DL) and limit of quantitation (LOQ).

APPL's laboratory control limits generated in house statistically do not meet the control limits listed in DoD QSM 4.2 for all analytes. Laboratory control spike recoveries for this project meet all control limits listed in the DoD QSM 4.2 except where noted. A copy of our in house generated control limits is available upon request. In addition, a copy of our LOQ control limits, established using 7 data points, are also available upon request.

Only the portion of the injection log relative to these samples is included. A full sequence log is available upon request.

Measurement uncertainty can be reported upon request.

The MADEP-EPH and VPH analyses were subcontracted to Gulf coast Analytical Laboratories, Inc. Their report is included.

CASE NARRATIVE

EPA Method 8015B

Total Petroleum Hydrocarbons – Diesel

Sample Preparation:

The water samples were extracted according to EPA method 3510C. The samples were extracted within holding time.

Sample Analysis Information:

The samples were analyzed according to the method using a Hewlett Packard Gas Chromatograph with a flame ionization detector.

Quality Control/Assurance

Calibrations:

Initial and continuing calibrations were performed according to the method. All calibration criteria were met.

Blanks:

No target analyte was detected above the detection limit in the method blank.

Spikes:

A Laboratory Control Spike (LCS) was used for quality assurance. All acceptance criteria were met.

Sample ES047 was designated by the client for MS/MSD analysis. Diesel fuel recovers below the 61% lower recovery limit in the MS at 33.5% and in the MSD at 30.0%.

Surrogates:

The surrogate recoveries are summarized on the form 2 & 8. All surrogate recoveries were within control limits.

Summary:

No other problem was encountered

EPA Method 8270D SIM

Polynuclear Aromatic Hydrocarbons

Sample Preparation:

The water samples were extracted according to EPA method 3510C. All holding times were met.

Sample Analysis Information:

The samples were analyzed according to EPA method 8270D using a Hewlett Packard Gas Chromatograph with a mass spectrometer detector in selective ion monitoring mode.

Quality Control/Assurance

Calibrations:

Initial and continuing calibrations were performed according to the method. All calibration criteria were met.

Blanks:

No target analyte was detected above the detection limits in the method blank.

Spikes:

A Laboratory Control Spike (LCS) was used for quality assurance. All spike criteria were met.

Sample ES047 was designated by the client for MS/MSD analysis. For the MSD, Acenaphthene recovered below the 45% lower control limit at 42.0%, Acenaphthylene below the 50% lower control limit at 45.6%, and Anthracene below the 55% lower control limit at 50.5%. The RPD for Naphthalene exceeded the 25% limit at 36.4%. All other recoveries met acceptance criteria.

Surrogates

Surrogate recoveries are summarized on the forms 2&8. All surrogate recoveries were within control limits.

Tuning:

The instrument was tuned using DFTPP. All method criteria were met.

Internal Standards

The internal standard area counts were compared to the mid-point of the initial calibration according to method 8270. All method criteria were met.

Summary:

No other problem was encountered.

EPA Method 8260B
Volatile Organic Analysis

Sample Preparation:

The water samples were purged according to EPA method 5030B. All holding times were met.

Sample Analysis Information:

The samples were analyzed according to the method using a Hewlett Packard Gas Chromatograph with a mass spectrometer detector. All samples were listed as non-preserved; they were injected within the seven day holding time. All holding times were met. Manual integrations were performed in accordance to APPL's SOP. All points of the gasoline curve, the gasoline second-source, and the gasoline LCS required manual integrations because the integration did not follow the baseline. Chromatograms of prior to and after manual integrations are enclosed.

Quality Control/Assurance:

Calibrations:

Initial and continuing calibrations were performed according to the method. All system performance check compounds and calibration check compounds met DoD acceptance criteria. All calibration criteria were met except for acetone in continuing calibration 1030C27W.D; acetone recovered above the 20% Drift limit at 28% Drift. Acetone recovered within the control limits of the following LCS and was not detected in any sample.

Blanks:

No target analyte was detected above the detection limits in the method blanks.

Spikes:

A Laboratory Control Spike (LCS) was used for quality assurance. A second-source standard was used for the LCS. All LCS recoveries were acceptable.

Sample ES047 was designated by the client for MS/MSD analysis. All spike criteria were met except for gasoline, which recovered above the 125% upper control limit at 126% in the MSD.

Surrogates:

Surrogate recoveries are summarized on Form 2 & 8. All surrogate recoveries were within the acceptance limits.

Tuning:

The instrument was tuned using BFB. All method criteria were met.

Internal Standards:

The internal standard area counts were compared to the mid-point of the initial calibration according to method 8260. All method criteria were met.

Summary:

No other problem was encountered. The data generated are acceptable.

EPA Method 6020
Dissolved Lead

Digestion Information:

The water samples were digested according to EPA methods 3015. All holding times were met.

Analysis Information:

Samples:

The samples were analyzed for dissolved lead according to EPA method 6020 using an Agilent 7500CX ICP-MS.

Calibrations:

The initial and continuing calibrations were analyzed according to the DoD QSM. The initial calibration verification is prepared from a second source standard.

Blanks:

No metal was detected at or above one-half the LOQ in the method blank.

Spikes:

Laboratory Control Spike (LCS), matrix spikes (MS/MSD), post digestion spike (PDS), and serial dilution were used for quality assurance. All LCS recoveries were within the acceptance limits.

Sample ES047 was designated by the client for MS/MSD analysis. All acceptance criteria were met in the MS/MSD, PDS, and DT.

Summary:

No analytical exception is noted.

Abbreviations and Flags

FLAG	DESCRIPTION
#	Recovery or RPD outside control limits
*	Recovery or RPD outside control limits
B	Analyte detected in associated method blank
C1	Reason for correction: wrote incorrect response
C2	Reason for correction: calculated incorrectly
C3	Reason for correction: needs to be rechecked
C4	Reason for correction: data not usable
DO	Diluted out
E	Exceeds linear range
F	Estimated value
G1	Includes a wide range of hydrocarbons which does not match our gasoline standard
G10	Includes a match to hydrocarbon profiles within the range of mineral spirits
G11	Includes a match to hydrocarbon profiles within the range of JP-4
G12	Pattern does not match the gasoline standard; the carbon range for this sample is consistent with JP8
G13	Closely resembles the hydrocarbon profile of aviation gasoline
G14	Analyte concentration may be biased due to carry over
G2	Closely resembles the boiling point hydrocarbon profile consistent with weathered gasoline
G3	Includes higher boiling hydrocarbons
G4	Includes dominant peak(s) not indicative of petroleum hydrocarbons
G5	Is mainly dominant peak(s) not indicative of petroleum hydrocarbons
G6	Contains recognizable contaminant peak(s) which has been removed from quantitation
G7	Is mainly a match to hydrocarbons within the range of gasoline
G8	Closely resembles the boiling point hydrocarbon profile consistent with weathered gasoline
G9	Includes hydrocarbons within the range of kerosene
J	Estimated value
M	Matrix effect
M1	Manual integration: integration does not follow baseline
M2	Manual integration: non-target peak interference
M3	Manual integration: to split a peak that was integrated as one peak by the computer
M4	Manual integration: to integrate a split peak
M5	Manual integration: the whole peak or part of the peak was not integrated
M6	Manual integration: computer integrated wrong peak
M7	Manual integration: other - explain
MDL	Method detection limit
ND	Not detected
NT	Non-target
Q	Acceptance criteria not met
T1 I	Includes wide range of hydrocarbons not indicative of diesel
T1 M	Is mainly wide range of hydrocarbons not necessarily indicative of diesel
T2 I	Includes lower boiling hydrocarbons, i.e. mineral spirits, kerosene, stoddard solvent, white gas
T2 M	Is mainly lower boiling hydrocarbons, i.e. mineral spirits, kerosene, stoddard solvent, white gas
T3 I	Includes higher boiling hydrocarbons, i.e. asphaltene, waster oil, motor oil, or weathered diesel fuel
T3 M	Is mainly higher boiling hydrocarbons, i.e. asphaltene, waster oil, motor oil, or weathered diesel fuel
T4 I	Includes dominant peak(s) not indicative of hydrocarbons
T4 M	Is mainly dominant peak(s) not indicative of hydrocarbons
T5	Contains recognizable contaminant peak(s) which has been removed from quantitation
T6	Is mainly a match to hydrocarbons within range of diesel fuel
T7	Closely resembles the boiling point hydrocarbon profile consistent with diesel fuel
T8	Includes a match to hydrocarbon profiles within range of diesel and kerosene fuel
T9 I	Includes non-diesel hydrocarbons within boiling point range of diesel fuel
T9 M	Is mainly non-diesel hydrocarbons within boiling point range of diesel fuel.
Y	Percent difference between primary and confirmation column > 40%

CHAIN OF CUSTODY,
ARF, CRF, AND
CLIENT COMMUNICATION

APPL - Analysis Request Form 66102

Client: Environet, Inc.
Address: 650 Iwilei Rd, #204
Honolulu, HI 96817
Attn: Stacy Fineran
Phone: 808-833-2225 Fax: 808-833-2231
Job: RED HILL/1022-024
PO #: 1022-015
Chain of Custody (Y/N): Y # 35457
RAD Screen (Y/N): Y pH (Y/N): Y
Turn Around Type: 2 WEEKS

Received by: TBV
Date Received: 10/26/11 Time: 10:00
Delivered by: FED EX
Shuttle Custody Seals (Y/N): N Time Zone: HAST
Chest Temp(s): SEE CRF 2.5-3.0°C
Color: VOA,M-PURPINK,Q-ORYE
Samples Chilled until Placed in Refrig/Freezer: Y
Project Manager: Cynthia Clark
QC Report Type: DVP4/ADRDOD/HI
Due Date: 11/09/11

Comments:
14 day TAT for Form 1s & 30 day TAT for full package. VDupra@environetInc.com
1 pdf on CD or FTP (no hard copy), NO hard copy to LDC per Stacy 2-24-11
Guidance: DOD QSM, EDD: Exce & ADR
DoD Forms, J flag to DL, U flag at LOD
EDD ADR A1/A3 (ADR 8.3a unchecked) to VDupra@ & sfineran@environetinc.com
metals 6020: report Lead with 0.5ug/L RL
TPH-Diesel only; VOCs: include gasoline by 8260B
MA-EPH subcontracted to Gulf Coast Analytical.

Sample Distribution:
GC: 3-$SIMHC12W, 3-$TPETD2
Extractions: 3-_SEP004S, 3-_SEP011
VOA: 4-$86RHBF
Metals: 3-$602D(Pb)
Other: 3-_M3015, 3-SUB

Charges:

Invoice To:
same

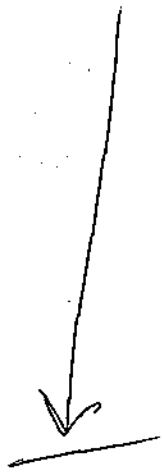
	Client ID		APPL ID	Sampled	Analyses Requested
1.	ES046		AY49333W	10/24/11 09:55	$602D(Pb), $86RHBF, $SIMHC12W, $TPETD2, SUB -- UN-PRESERVED VOA VIALS
2.	ES047	MS/MSD,LTDVOL	AY49334W	10/24/11 08:30	$602D(Pb), $86RHBF, $SIMHC12W, $TPETD2, SUB -- UN-PRESERVED VOA VIALS,LTD VOL on amber
3.	ES048		AY49335W	10/24/11 07:00	$86RHBF -- UN-PRESERVED VOA VIALS
4.	ES049		AY49336W	10/24/11 14:35	$602D(Pb), $86RHBF, $SIMHC12W, $TPETD2, SUB -- UN-PRESERVED VOA VIALS

APPL Sample Receipt Form

ARF# **66102**

Sample	Container Type	Count	pH
AY49333	[6] PL 500mL - HNO3	1	1.7
	[13] VOAs - HCL	3	na
	[15] VOAs - NP	4	na
	[17] Amber Liter	3	na
	[26] Other	2	1.7
AY49334	[6] PL 500mL - HNO3	4	1.7
	[13] VOAs - HCL	12	na
	[15] VOAs - NP	16	na
	[17] Amber Liter	9	na
	[26] Other	7	1.7
AY49335	[15] VOAs - NP	1	na
AY49336	[6] PL 500mL - HNO3	1	1.7
	[13] VOAs - HCL	3	na
	[15] VOAs - NP	4	na
	[17] Amber Liter	3	na
	[26] Other	2	1.7

Sample	Container Type	Count	pH
other - 1 Liter amber HCl			



66102
3.0

APPL LABS

APPL, Inc.
908 N Temperance Ave
Clovis, CA 93611

Phone: (559) 275-2175
Fax: (559) 275-4422

CHAIN OF CUSTODY RECORD

C.O.C. 35457

Report to: PLEASE PRINT		Invoice to: PLEASE PRINT	
Company Name: Environet	Phone: 808 833-2225	Company Name: Environet	Phone: 808-833-2225
Address: 650 Iwilei Rd. #204 Honolulu, Hi	Fax: 808 833-2231	Address: 650 Iwilei Rd. #204 Honolulu, Hi	Fax: 808 833-2231
Attn: Stacey Fineran		Attn: A.P. Trisha Yosuda	

Project Name/Number	Sampler (Print)	Purchase Order Number	Sampler (Signature)
Red Hill /1022-024	Stacey Fineran		

Sample Identification	Location	Date Collected	Time Collected	No. of Containers	Matrix: Aq	Sed.	Soil	Analysis Requested/Method Number: EPA 8015 B TPH DRO	EPA 8260 TPH-GRO	EPA 8260 VOC	EPA 8270C PAH	EPA 6020 Dissolved Lead	MADEP-EPH	MADEP-VPH
ES046	RHSF	10/24/11	0955	13	X			X	X	X	X	X	X	X
ES047MSMSD			0830	52	X			X	X	X	X	X	X	X
ES048			0700	1	X					X				
ES049	RHSF	10/24/11		13	X			X	X	X	X	X	X	X

Date Shipped: 10/25/2011
Carrier: Fed Express
Waybill No.:
Comments: *HCl was not rinsed out of VOAs

15

Shuttle Temperature:

Turnaround Requested: MUST CHECK ONE
☒ Standard (2-3 week) ☐ One week ☐ 24-48 hour

Sample Disposal:
☐ Return to client ☒ Disposal by Lab (30-day retention)

Relinquished by sampler:	Date	Time	Received by:	Relinquished by:	Date	Time	Received by:
	10/25	1400	Fed Ex				
Relinquished by:	Date	Time	Received by:	Relinquished by:	Date	Time	Received at lab by:
					10/26/11	1000	

White: Return to client with report Yellow: Laboratory Copy Pink: Sampler

See reverse side for Container Preservative and Sampling Information

COOLER RECEIPT FORM

1) Project: RED HILL / 1022-024 Date Received: 10/26/11

2) Coolers: Number of Coolers: 3

3) YES NO Were coolers and samples screened for radioactivity?

4) YES NO Were custody seals on outside of cooler? How many? _____ Date on seal? _____

5) Name on seal? _____

6) YES NO NA Were custody seals unbroken and intact at the time of arrival?

7) YES NO Did the cooler come with a shipping slip (air bill, etc.)? Carrier name: Fed Ex

8) Shipping slip numbers: 1) 8768 4695 9176 (MASTER) 2) _____ 3) _____

9) YES NO NA Was the shipping slip scanned into the database?

10) YES NO NA If cooler belongs to APPL, has it been logged into the ice chest database?

11) Describe type of packing in cooler (bubble wrap, popcorn, type of ice, etc.): bubble bag in wet ice

12) YES NO NA For hand delivered samples was sufficient ice present to start the cooling process?

13) YES NO Was a temperature blank included in the cooler?

14) Serial number of certified NIST thermometer used: A39267 Correction factor: 0

15) Cooler temp(s): 1) 2.5°C 2) 3.0°C 3) 3.0°C 4) _____ 5) _____ 6) _____ 7) _____ 8) _____

Chain of custody:

16) YES NO Was a chain of custody received?

17) YES NO Were the custody papers signed in the appropriate places?

18) YES NO Was the project identifiable from custody papers?

19) YES NO Did the chain of custody include date and time of sampling?

20) YES NO Is location where sample was taken listed on the chain of custody?

Sample Labels:

21) YES NO Were container labels in good condition?

22) YES NO Was the client ID on the label?

23) YES NO Was the date of sampling on the label?

24) YES NO Was the time of sampling on the label?

25) YES NO Did all container labels agree with custody papers?

Sample Containers:

26) YES NO Were all containers sealed in separate bags?

27) YES NO Did all containers arrive unbroken?

28) YES NO Was there any leakage from samples?

29) YES NO Were any of the lids cracked or broken?

30) YES NO Were correct containers used for the tests indicated?

31) YES NO Was a sufficient amount of sample sent for tests indicated?

32) YES NO NA Were bubbles present in volatile samples? If yes, the following were received with air bubbles:

Larger than a pea: YZ 10/26/11

Smaller than a pea: AY49333W03, ~~AY49334W30~~ AY49334W26-W29, AY49335W01

Preservation & Hold time:

33) YES NO NA Was a sufficient amount of holding time remaining to analyze the samples?

34) YES NO NA Do the sample containers contain the same preservative as what is stated on the COC?

35) YES NO NA Was the pH taken of all non-VOA preserved samples and written on the sample container?

36) YES NO NA Was the pH of acid preserved non-VOA samples < 2 & sodium hydroxide preserved samples > 12?

37) YES NO NA Unpreserved VOA Vials received? _____

38) YES NO NA Are unpreserved VOA vials noted in the ADD TEST FIELD on the ARF? _____

Lab notified if pH was not adequate: _____

Deficiencies: Received 4 Amber liter broken, 9 left for Sample ES047 ms/msd

Signature of personnel receiving samples: _____ Second reviewer: _____

Signature of project manager notified: Renée Date and Time of notification: 10-26-11

Name of client notified: _____ Date and Time of notification: _____

Information given to client: _____

_____ by whom (initials): _____

EPA 8015 Modified
Total Petroleum Hydrocarbons

EPA 8015 Modified
Total Petroleum Hydrocarbons
QC Summary

Method Blank
TPH Diesel Water

Blank Name/QCG: **111031W-49334 - 160886**
Batch ID: #TPETD-111031A

APPL Inc.
908 North Temperance Avenu
Clovis, CA 93611

Sample Type	Analyte	Result	LOQ	LOD	DL	Units	Extraction Date	Analysis Date
BLANK	DIESEL FUEL	80.8 U	150	80.8	40.4	ug/L	10/31/11	11/06/11
BLANK	SURROGATE: OCTACOSANE (S)	71.2	28-142			%	10/31/11	11/06/11
BLANK	SURROGATE: ORTHO-TERPHEN	60.5	57-132			%	10/31/11	11/06/11

Quant Method: TPH1028.M
Run #: 1106005
Instrument: Apollo
Sequence: 111106
Initials: LA

GC SC-Blank-REG MDLs
Printed: 11/30/11 11:27:27 AM

Form 2 & 8

Surrogate Recovery

Lab Name: APPL, Inc. SDG No: 66102

Case No: 66102 Date Analyzed: 11/06/11

Matrix: WATER Instrument: Apollo

APPL ID.	Client Sample No.	SURROGATE: OCTACOSANE (S)			SURROGATE: ORTHO-TERPHENYL (S)		
		Limits	Result	Qualifier	Limits	Result	Qualifier
111031A-BLK	Blank	28-142	71.2		57-132	60.5	
AY49333	ES046	28-142	97.0		57-132	69.1	
AY49334-MS	Matrix Spike	28-142	87.3		57-132	72.0	
AY49334-MSD	Matrix SpikeD	28-142	83.3		57-132	59.5	
AY49334	ES047	28-142	107		57-132	72.2	
AY49336	ES049	28-142	106		57-132	63.4	
111031A-LCS	Lab Control Spike	28-142	83.3		57-132	98.0	

Comments: Batch: #TPETD-111031A

Laboratory Control Spike Recovery
TPH Diesel Water

APPL ID: **111031W-49334 LCS - 160886**
Batch ID: #TPETD-111031A

APPL Inc.
908 North Temperance Avenue
Clovis, CA 93611

Compound Name	Spike Level ug/L	SPK Result ug/L	SPK % Recovery	Recovery Limits
DIESEL FUEL	2000	1520	76.0	61-143
SURROGATE: OCTACOSANE (S)	150	125	83.3	28-142
SURROGATE: ORTHO-TERPHENYL (S)	150	147	98.0	57-132

Comments:

Primary	SPK
Quant Method :	TPH8S15.M
Extraction Date :	10/31/11
Analysis Date :	11/29/11
Instrument :	Apollo
Run :	1129017
Initials :	LA

Printed: 11/30/11 11:27:20 AM

APPL Standard LCS

Matrix Spike Recoveries
TPH Diesel Water

APPL ID: **111031W-49334 MS - 160886**
Batch ID: #TPETD-111031A
Sample ID: AY49334
Client ID: ES047

APPL Inc.
908 North Temperance Avenue
Clovis, CA 93611

Compound Name	Spike Lvl ug/L	Matrix Result ug/L	SPK Result ug/L	DUP Result ug/L	SPK % Recovery	DUP % Recovery	Recovery Limits	RPD %	RPD Limits
DIESEL FUEL	2000	730	1400	1330	33.5 #	30.0 #	61-143	5.1	30
SURROGATE: OCTACOSANE (S)	150	NA	131	125	87.3	83.3	28-142		
SURROGATE: ORTHO-TERPHENYL (S)	150	NA	108	89.3	72.0	59.5	57-132		

= Recovery is outside QC limits.

Comments:

Primary	SPK	DUP
Quant Method :	TPH1028.M	TPH1028.M
Extraction Date :	10/31/11	10/31/11
Analysis Date :	11/06/11	11/06/11
Instrument :	Apollo	Apollo
Run :	1106014	1106015
Initials :	LA	

Printed: 11/30/11 11:27:17 AM
APPL MSD SCII

EPA 8015B-e

Form 4

Blank Summary

Lab Name: APPL, Inc. SDG No: 66102

Case No: 66102 Date Analyzed: 11/06/11

Matrix: WATER Instrument: Apollo

Blank ID: 111031A-BLK Time Analyzed: 1722

APPL ID.	Client Sample No.	File ID.	Date Analyzed
111031A-BLK	Blank	1106005	11/06/11 1722
AY49333	ES046	1106013	11/06/11 2030
111031A-MS	Matrix Spike	1106014	11/06/11 2053
111031A-MSD	Matrix SpikeD	1106015	11/06/11 2117
AY49334	ES047	1106016	11/06/11 2140
AY49336	ES049	1106019	11/06/11 2250
111031A-LCS	Lab Control Spike	1129017	11/29/11 1845

Comments: Batch: #TPETD-111031A

EPA 8015 Modified
Total Petroleum Hydrocarbons
Sample Data

TPH Diesel Water

Environet, Inc.
650 Iwilei Rd, #204
Honolulu, HI 96817

Attn: Stacy Fineran
Project: RED HILL/1022-024
Sample ID: ES046
Sample Collection Date: 10/24/11

APPL Inc.
908 North Temperance Avenue
Clovis, CA 93611

ARF: 66102
APPL ID: AY49333
QCG: #TPETD-111031A-160886

Method	Analyte	Result	LOQ	LOD	DL	Units	Extraction Date	Analysis Date
EPA 8015B-	DIESEL FUEL	750 ++	150	80.8	40.4	ug/L	10/31/11	11/06/11
EPA 8015B-	SURROGATE: OCTACOSANE (S)	97.0	28-142			%	10/31/11	11/06/11
EPA 8015B-	SURROGATE: ORTHO-TERPHENYL (S	69.1	57-132			%	10/31/11	11/06/11

++(T7) The analyst has noted that the chromatogram of this sample closely resembles the boiling point hydrocarbon profile consistent with diesel fuel.

Quant Method: TPH1028.M
Run #: 1106013
Instrument: Apollo
Sequence: 111106
Dilution Factor: 1
Initials: LA

Printed: 11/30/11 11:27:24 AM
APPL-F1-SC-NoMC-REG MDLs

Quantitation Report (QT Reviewed)

Data File : G:\APOLLO\DATA\111106\1106013.D Vial: 13
Acq On : 11-6-11 20:30:17 Operator: LAC
Sample : AY49333W09 5/1030 Inst : Apollo
Misc : Water Multiplr: 4.85
IntFile : events.e
Quant Time: Nov 7 9:50 2011 Quant Results File: TPH1028.RES

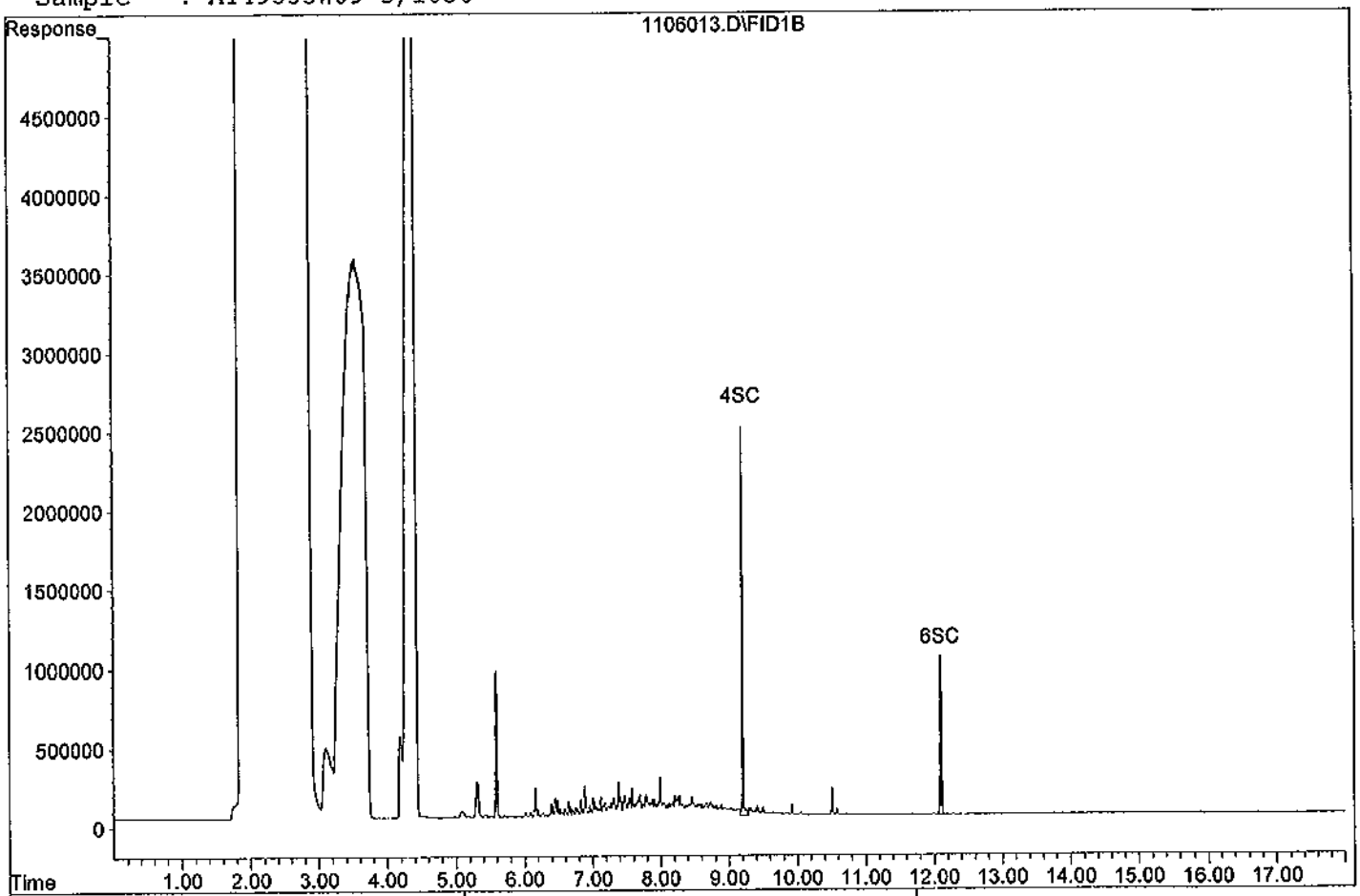
Method : G:\APOLLO\DATA\111028\TPH1028.M (Chemstation Integrator)
Title : Diesel
Last Update : Mon Oct 31 10:02:11 2011
Response via : Multiple Level Calibration

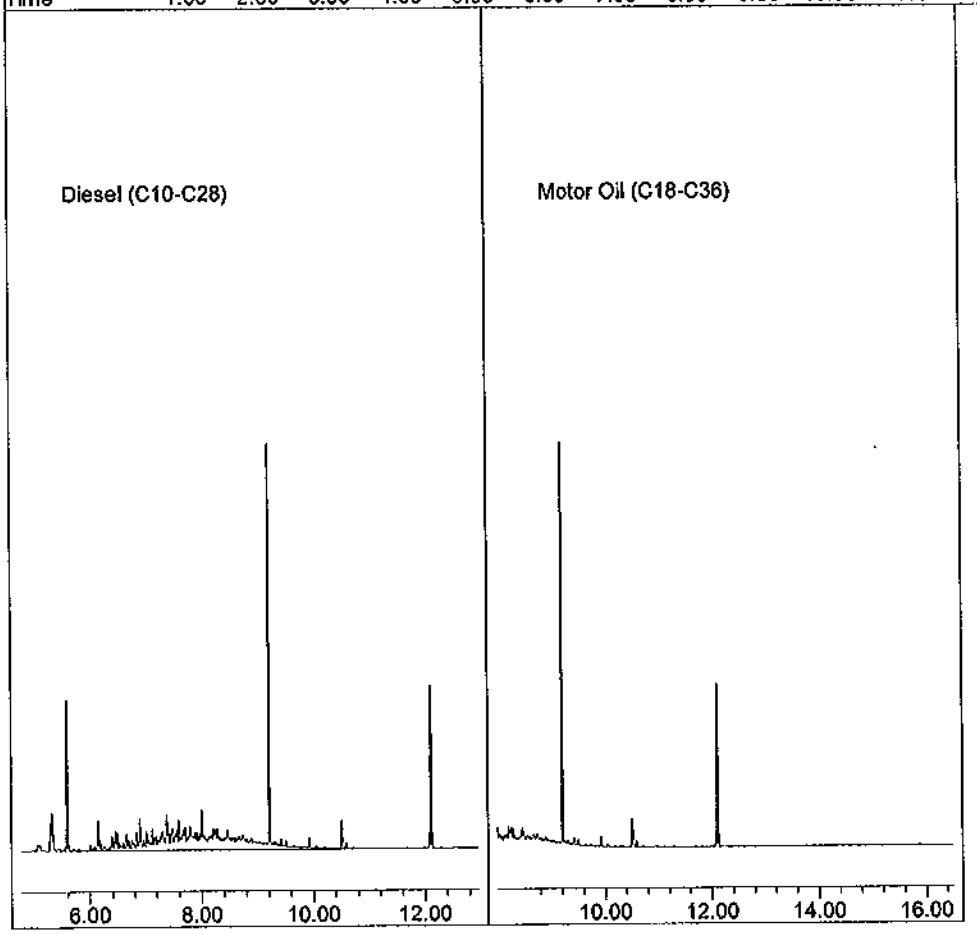
Volume Inj. : 2UL
Signal Phase : DB-5
Signal Info : FID02A

Compound	R.T.	Response	Conc Units
System Monitoring Compounds			
4) SC Ortho-Terphenyl(S)	9.21	18082118	100.664 ppb
Surrogate Spike 145.631		Recovery =	69.12%
6) SC Octacosane(S)	12.10	13534363	141.185 ppb
Surrogate Spike 145.631		Recovery =	96.95%
Target Compounds			
1) HATM Diesel (C10-C28)	8.86	130505484	752.499 ppb

TT LAC 11/30/11

Data File: G:\APOLLO\DATA\111106\1106013.D
Sample : AY49333W09 5/1030





TPH Diesel Water

Environet, Inc.
650 Iwilei Rd, #204
Honolulu, HI 96817

Attn: Stacy Fineran
Project: RED HILL/1022-024
Sample ID: ES047
Sample Collection Date: 10/24/11

APPL Inc.
908 North Temperance Avenue
Clovis, CA 93611

ARF: 66102
APPL ID: AY49334
QCG: #TPETD-111031A-160886

Method	Analyte	Result	LOQ	LOD	DL	Units	Extraction Date	Analysis Date
EPA 8015B-	DIESEL FUEL	730 ++	150	80.8	40.4	ug/L	10/31/11	11/06/11
EPA 8015B-	SURROGATE: OCTACOSANE (S)	107	28-142			%	10/31/11	11/06/11
EPA 8015B-	SURROGATE: ORTHO-TERPHENYL (S	72.2	57-132			%	10/31/11	11/06/11

++(T7) The analyst has noted that the chromatogram of this sample closely resembles the boiling point hydrocarbon profile consistent with diesel fuel.

Quant Method:	TPH1028.M
Run #:	1106016
Instrument:	Apollo
Sequence:	111106
Dilution Factor:	1
Initials:	LA

Printed: 11/30/11 11:27:24 AM
APPL-F1-SC-NoMC-REG MDLs

Quantitation Report (QT Reviewed)

Data File : G:\APOLLO\DATA\111106\1106016.D Vial: 16
Acq On : 11-6-11 21:40:27 Operator: LAC
Sample : AY49334W32 5/1030 Inst : Apollo
Misc : Water Multiplr: 4.85
IntFile : events.e
Quant Time: Nov 7 9:50 2011 Quant Results File: TPH1028.RES

Method : G:\APOLLO\DATA\111028\TPH1028.M (Chemstation Integrator)
Title : Diesel
Last Update : Mon Oct 31 10:02:11 2011
Response via : Multiple Level Calibration

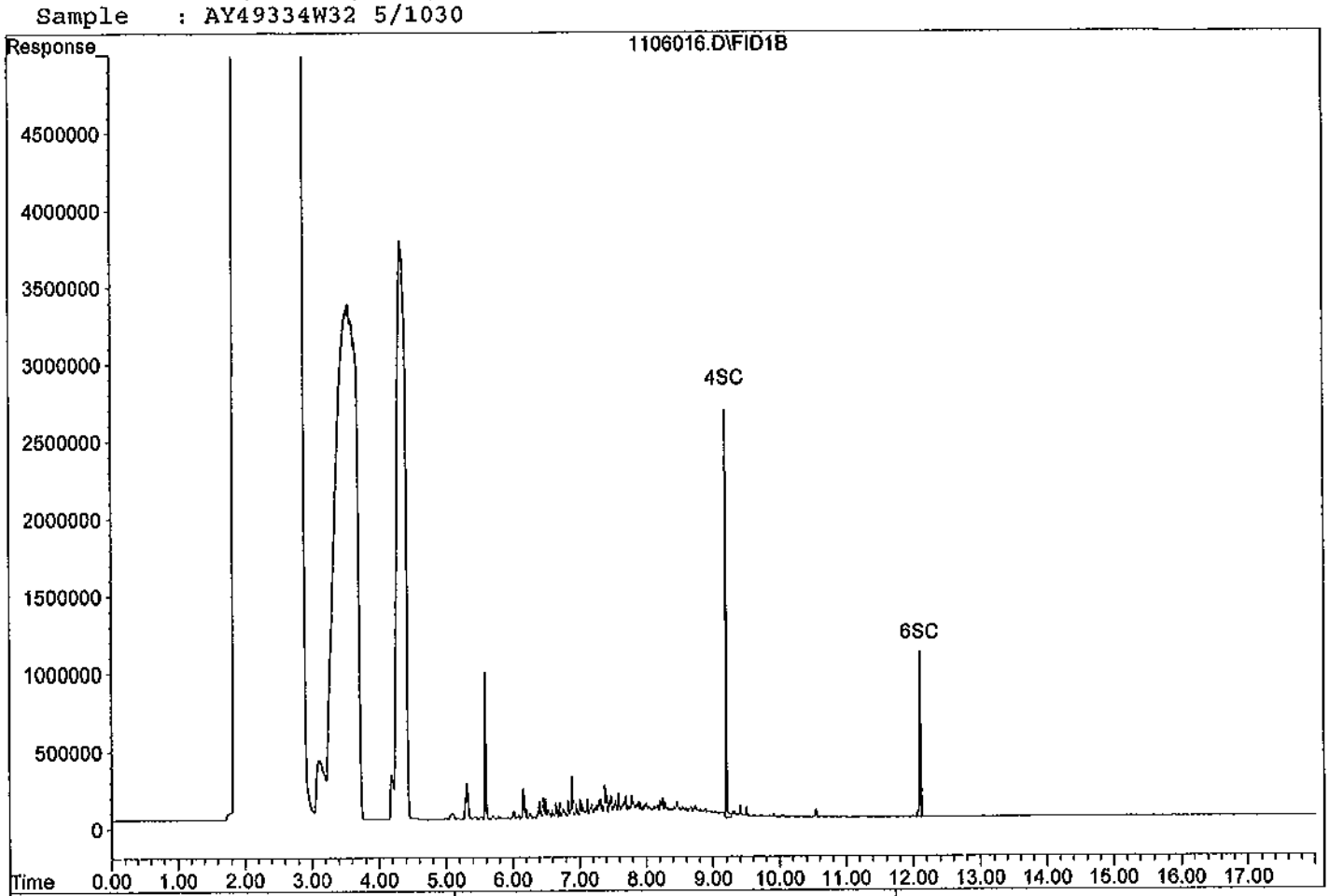
Volume Inj. : 2UL
Signal Phase : DB-5
Signal Info : FID02A

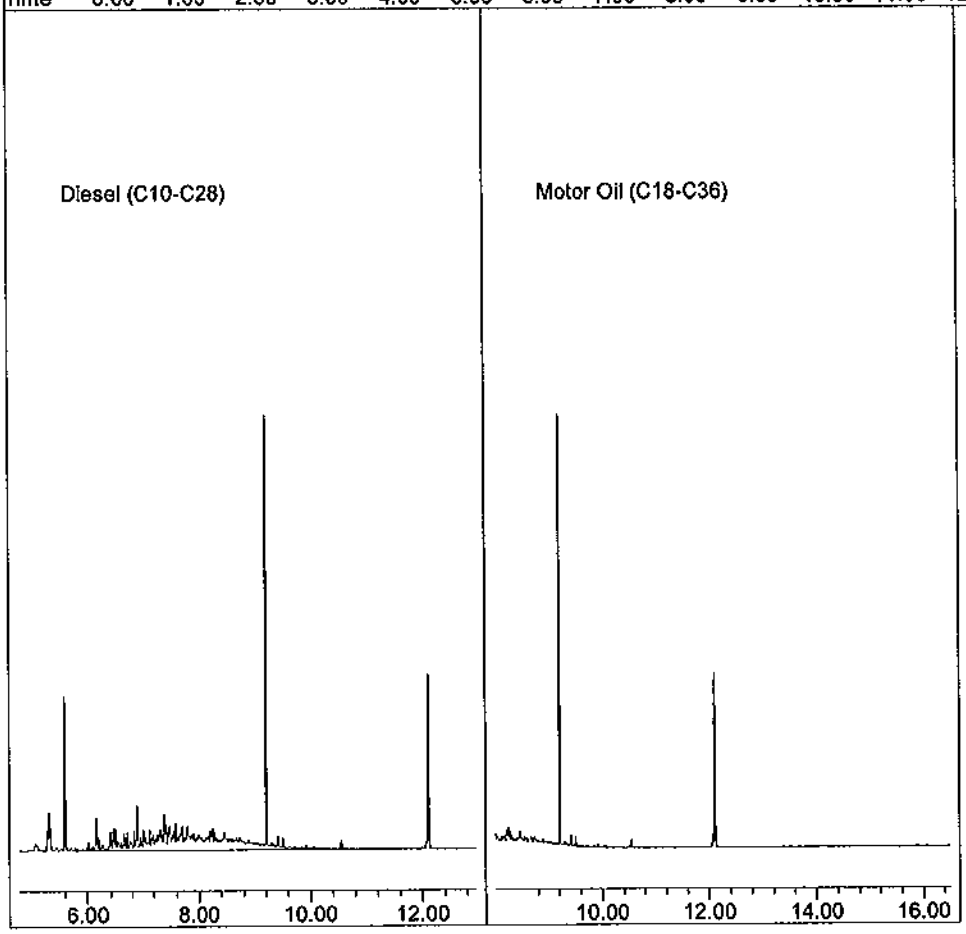
Compound	R.T.	Response	Conc Units
System Monitoring Compounds			
4) SC Ortho-Terphenyl(S)	9.21	18885119	105.134 ppb
Surrogate Spike 145.631		Recovery =	72.19%
6) SC Octacosane(S)	12.10	14910092	155.536 ppb
Surrogate Spike 145.631		Recovery =	106.80%
Target Compounds			
1) HATM Diesel (C10-C28)	8.86	125913483	726.021 ppb

T7 LAC 11/30/11

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

Data File: G:\APOLLO\DATA\111106\1106016.D
Sample : AY49334W32 5/1030





TPH Diesel Water

Environet, Inc.
650 Iwilei Rd, #204
Honolulu, HI 96817

Attn: Stacy Fineran
Project: RED HILL/1022-024
Sample ID: ES049
Sample Collection Date: 10/24/11

APPL Inc.
908 North Temperance Avenue
Clovis, CA 93611

ARF: 66102
APPL ID: AY49336
QCG: #TPETD-111031A-160886

Method	Analyte	Result	LOQ	LOD	DL	Units	Extraction Date	Analysis Date
EPA 8015B-	DIESEL FUEL	80.8 U	150	80.8	40.4	ug/L	10/31/11	11/06/11
EPA 8015B-	SURROGATE: OCTACOSANE (S)	106	28-142			%	10/31/11	11/06/11
EPA 8015B-	SURROGATE: ORTHO-TERPHENYL (S	63.4	57-132			%	10/31/11	11/06/11

Quant Method: TPH1028.M
Run #: 1106019
Instrument: Apollo
Sequence: 111106
Dilution Factor: 1
Initials: LA

Printed: 11/30/11 11:27:24 AM
APPL-F1-SC-NoMC-REG MDLs

Quantitation Report (QT Reviewed)

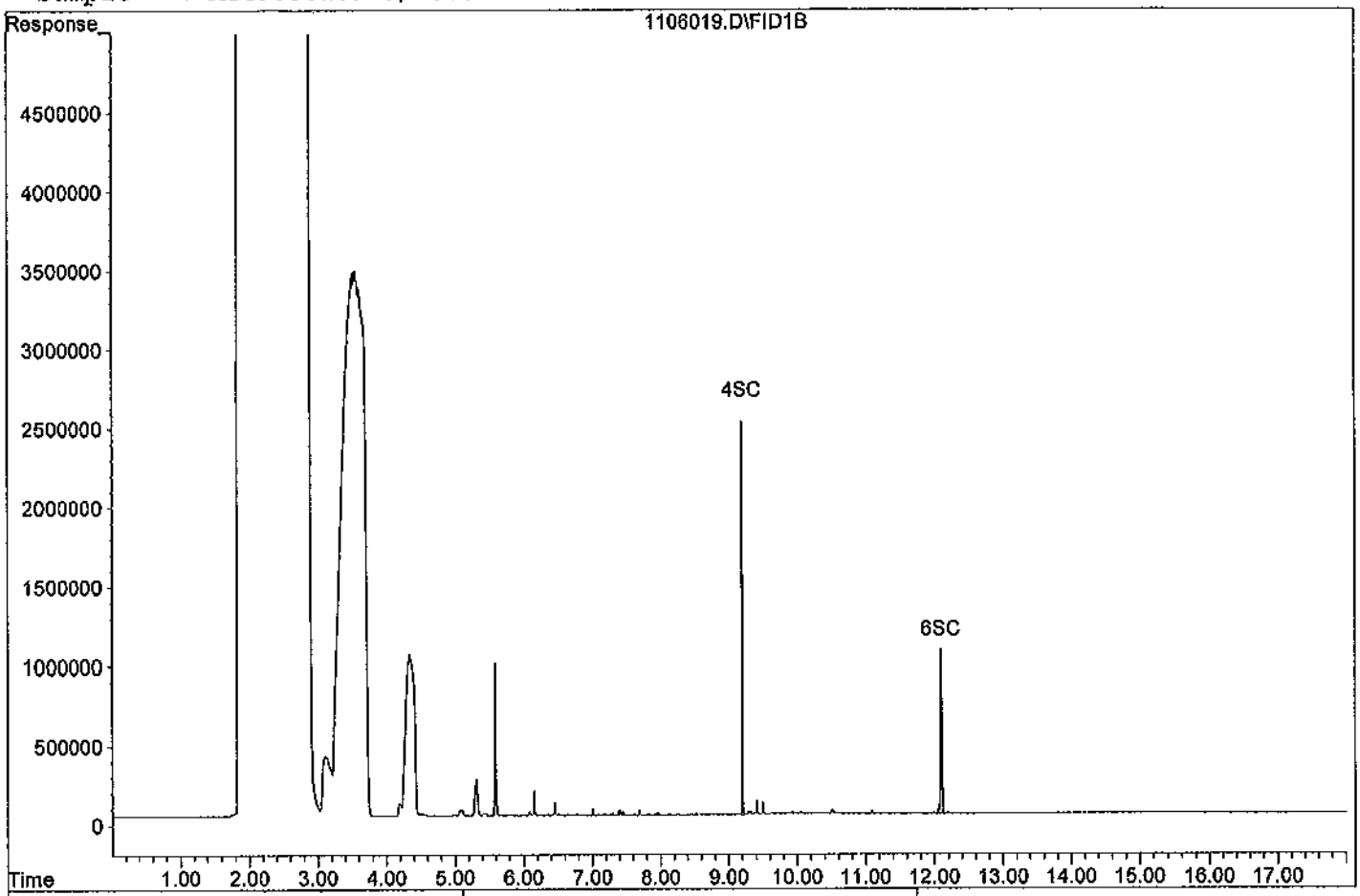
Data File : G:\APOLLO\DATA\111106\1106019.D Vial: 19
Acq On : 11-6-11 22:50:22 Operator: LAC
Sample : AY49336W09 5/1050 Inst : Apollo
Misc : Water Multiplr: 4.76
IntFile : events.e
Quant Time: Nov 7 10:06 2011 Quant Results File: TPH1028.RES

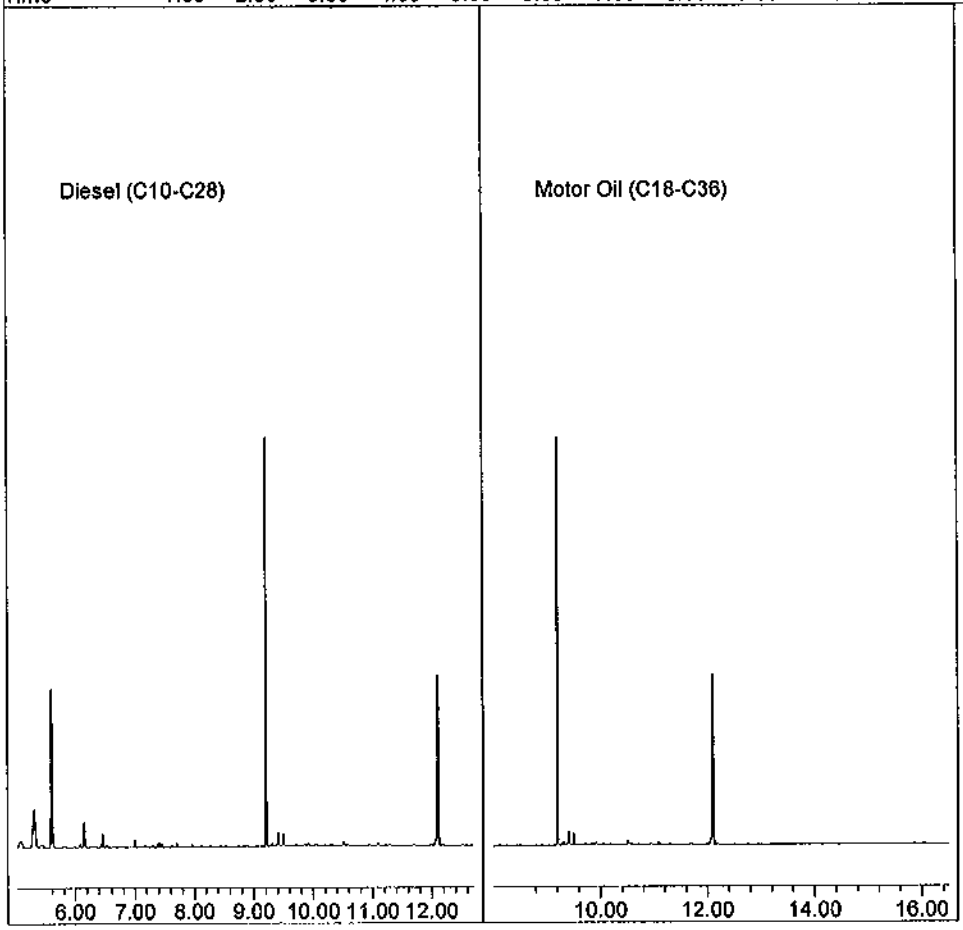
Method : G:\APOLLO\DATA\111028\TPH1028.M (Chemstation Integrator)
Title : Diesel
Last Update : Mon Oct 31 10:02:11 2011
Response via : Multiple Level Calibration

Volume Inj. : 2UL
Signal Phase : DB-5
Signal Info : FID02A

Compound	R.T.	Response	Conc Units
System Monitoring Compounds			
4) SC Ortho-Terphenyl(S)	9.20	16586813	90.580 ppb
Surrogate Spike 142.857		Recovery =	63.41%
6) SC Octacosane(S)	12.10	14736715	150.799 ppb
Surrogate Spike 142.857		Recovery =	105.56%
Target Compounds			

Data File: G:\APOLLO\DATA\111106\1106019.D
Sample : AY49336W09 5/1050





EPA 8015 Modified
Total Petroleum Hydrocarbons

Calibration Data

TPH Extractables

TPH1028

Form 6

Initial Calibration

Lab Name: APPL, Inc. SDG No: 66102

Case No: Initial Cal. Date: 10/28/11

Matrix: Instrument: Apollo Initials: LAC

			1028016.D	1028017.D	1028018.D	1028019.D	1028020.D	1028021.D								
		Compound	1	2	3	4	5	6					Avg	%RSD		
1	HATM	Diesel (C10-C28)	442719	420460	472596	423976	425504	340418					420946	10	HATM	
2	HBTM	Motor Oil (C18-C36)	178423	183716	183814	180648	189866	171166					181272	3.5	HBTM	
3	SA	Not Used(S)	492314	497998	549709	499503	505363	680489					537563	14	SA	
4	SC	Ortho-Terphenyl(S)		432362	446201	461765	440594	399046					435993	5.3	SC	
5	SA	Not Used2(S)	246114	255075	272188	252864	246503	213571					247719	7.8	SA	
6	SC	Octacosane(S)		229602	230817	244618	233443	224899					232676	3.2	SC	
7																
8																
9																
10																
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1.2501454

Quantitation Report (Not Reviewed)

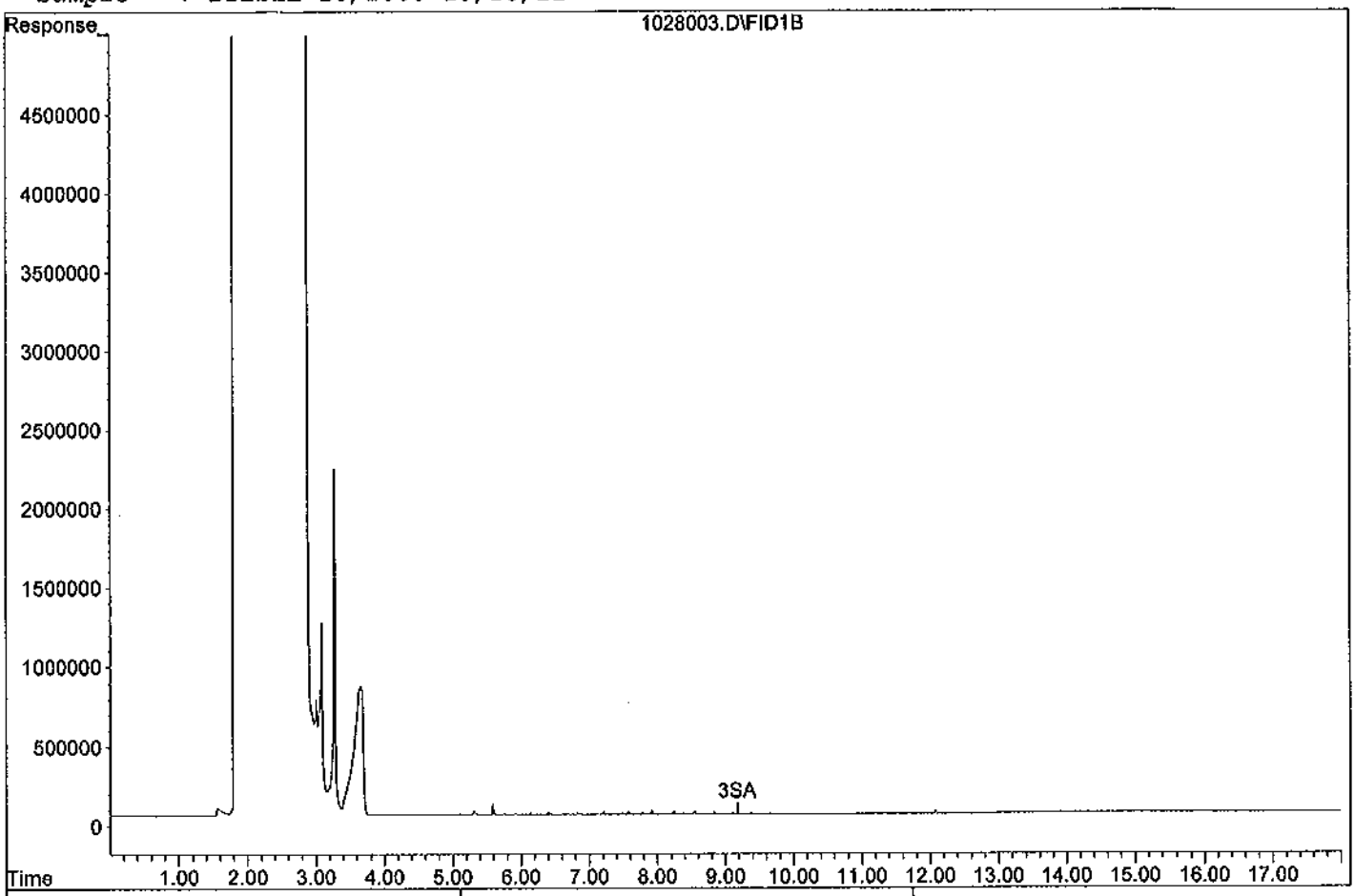
Data File : G:\APOLLO\DATA\111028\1028003.D Vial: 3
Acq On : 10-28-11 9:47:18 Operator: LAC
Sample : DIESEL 10/1000 10/28/11 Inst : Apollo
Misc : Mix(A) Multiplr: 1.00
IntFile : events.e
Quant Time: Oct 28 13:43 2011 Quant Results File: TPH1028.RES

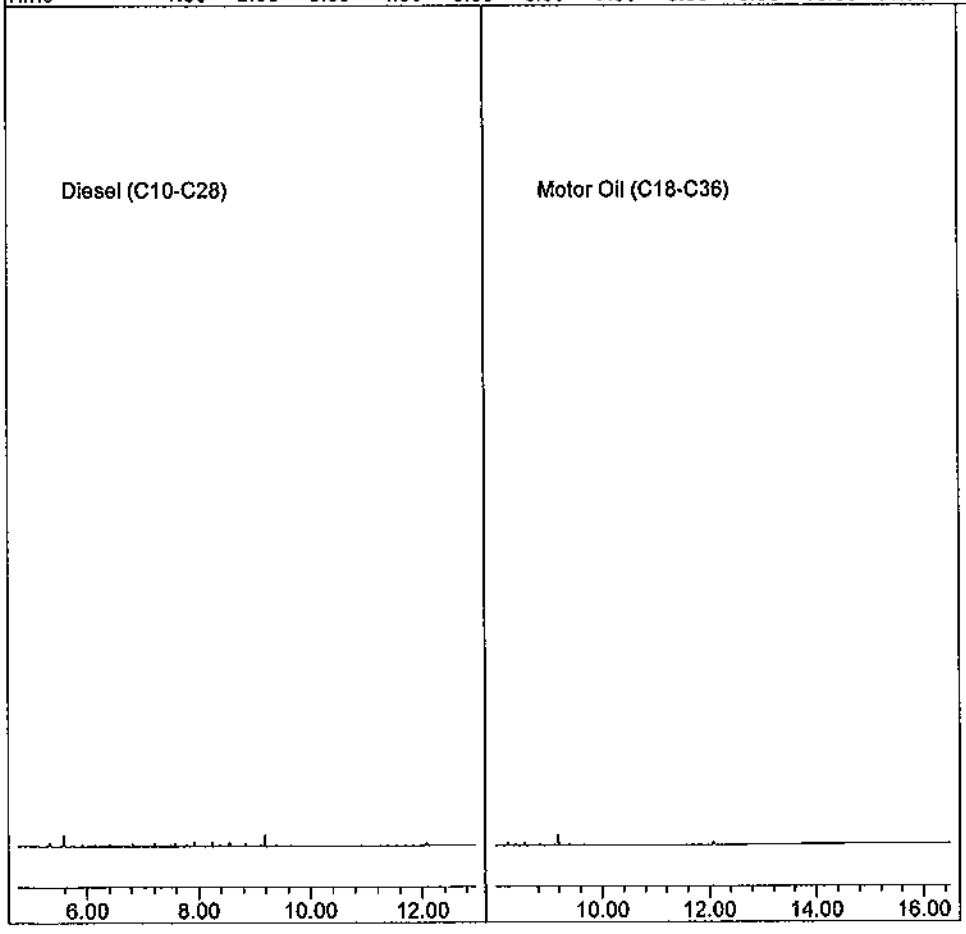
Method : G:\APOLLO\DATA\111028\TPH1028.M (Chemstation Integrator)
Title : Diesel
Last Update : Mon Oct 31 10:02:11 2011
Response via : Multiple Level Calibration

Volume Inj. : 2UL
Signal Phase : DB-5
Signal Info : FID02A

Compound	R.T.	Response	Conc Units
System Monitoring Compounds			
3) SA Not Used(S)	9.18	492314	0.777 ppb
Surrogate Spike 30.000	Recovery	=	2.59%
Target Compounds			
1) HATM Diesel (C10-C28)	8.86	8854372	10.215 ppb

Data File: G:\APOLLO\DATA\111028\1028003.D
Sample : DIESEL 10/1000 10/28/11





Quantitation Report (Not Reviewed)

Data File : G:\APOLLO\DATA\111028\1028004.D Vial: 4
Acq On : 10-28-11 10:11:19 Operator: LAC
Sample : DIESEL 100/1000 Inst : Apollo
Misc : Mix(A) Multiplr: 1.00
IntFile : events.e
Quant Time: Oct 28 13:43 2011 Quant Results File: TPH1028.RES

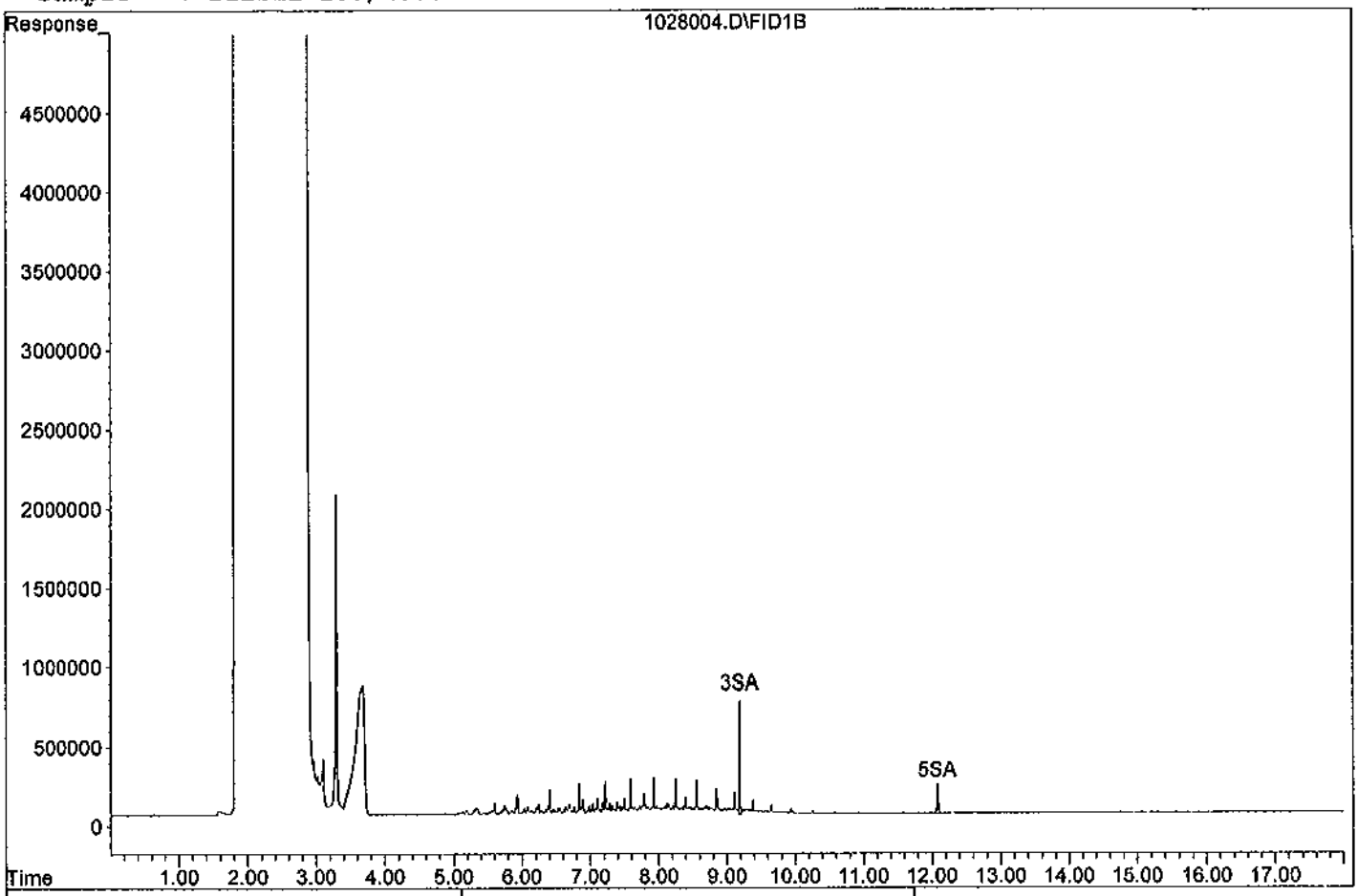
Method : G:\APOLLO\DATA\111028\TPH1028.M (Chemstation Integrator)
Title : Diesel
Last Update : Mon Oct 31 10:02:11 2011
Response via : Multiple Level Calibration

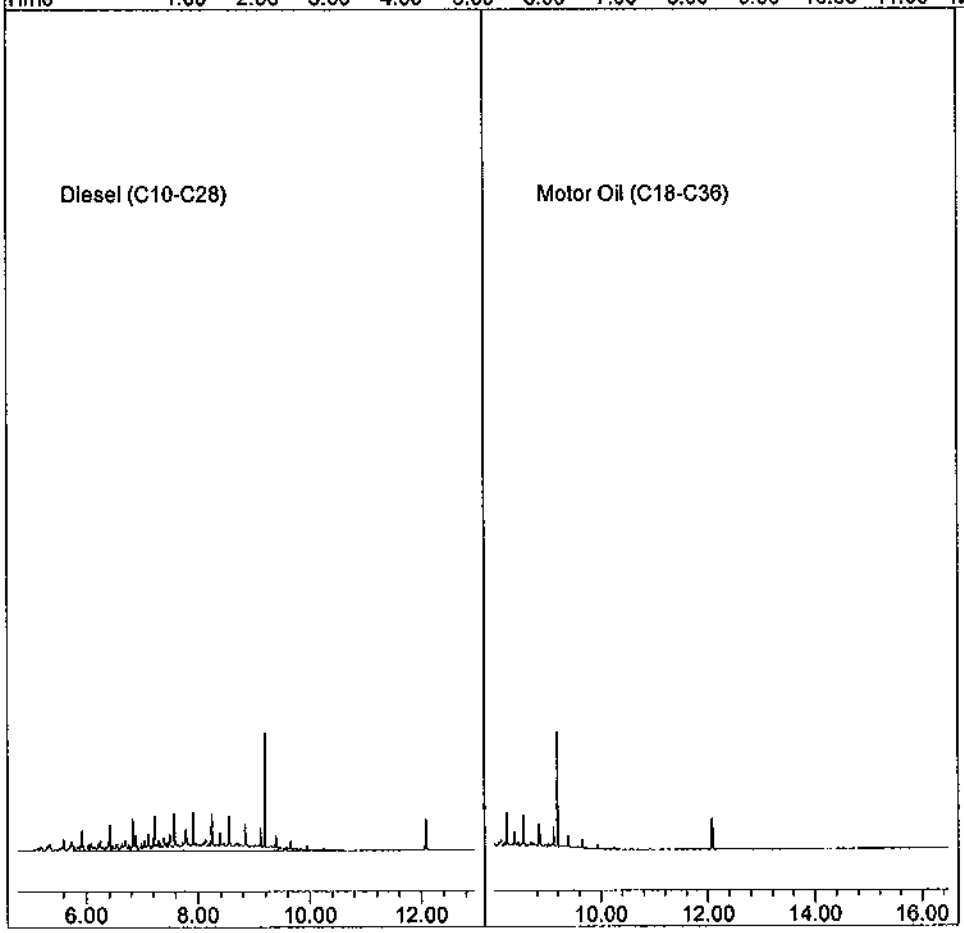
Volume Inj. : 2UL
Signal Phase : DB-5
Signal Info : FID02A

Compound	R.T.	Response	Conc Units
System Monitoring Compounds			
3) SA Not Used(S)	9.19	4979981	7.855 ppb
Surrogate Spike 30.000	Recovery =		26.18%
5) SA Not Used2(S)	12.08	2550752	9.065 ppb
Surrogate Spike 30.000	Recovery =		30.22%
Target Compounds			
1) HATM Diesel (C10-C28)	8.86	84092037	97.013 ppb

(f)=RT Delta > 1/2 Window 38 (m)=manual int.

Data File: G:\APOLLO\DATA\111028\1028004.D
Sample : DIESEL 100/1000





Quantitation Report (Not Reviewed)

Data File : G:\APOLLO\DATA\111028\1028005.D Vial: 5
Acq On : 10-28-11 10:35:26 Operator: LAC
Sample : DIESEL 400/1000 Inst : Apollo
Misc : Mix(A) Multiplr: 1.00
IntFile : events.e
Quant Time: Oct 28 13:43 2011 Quant Results File: TPH1028.RES

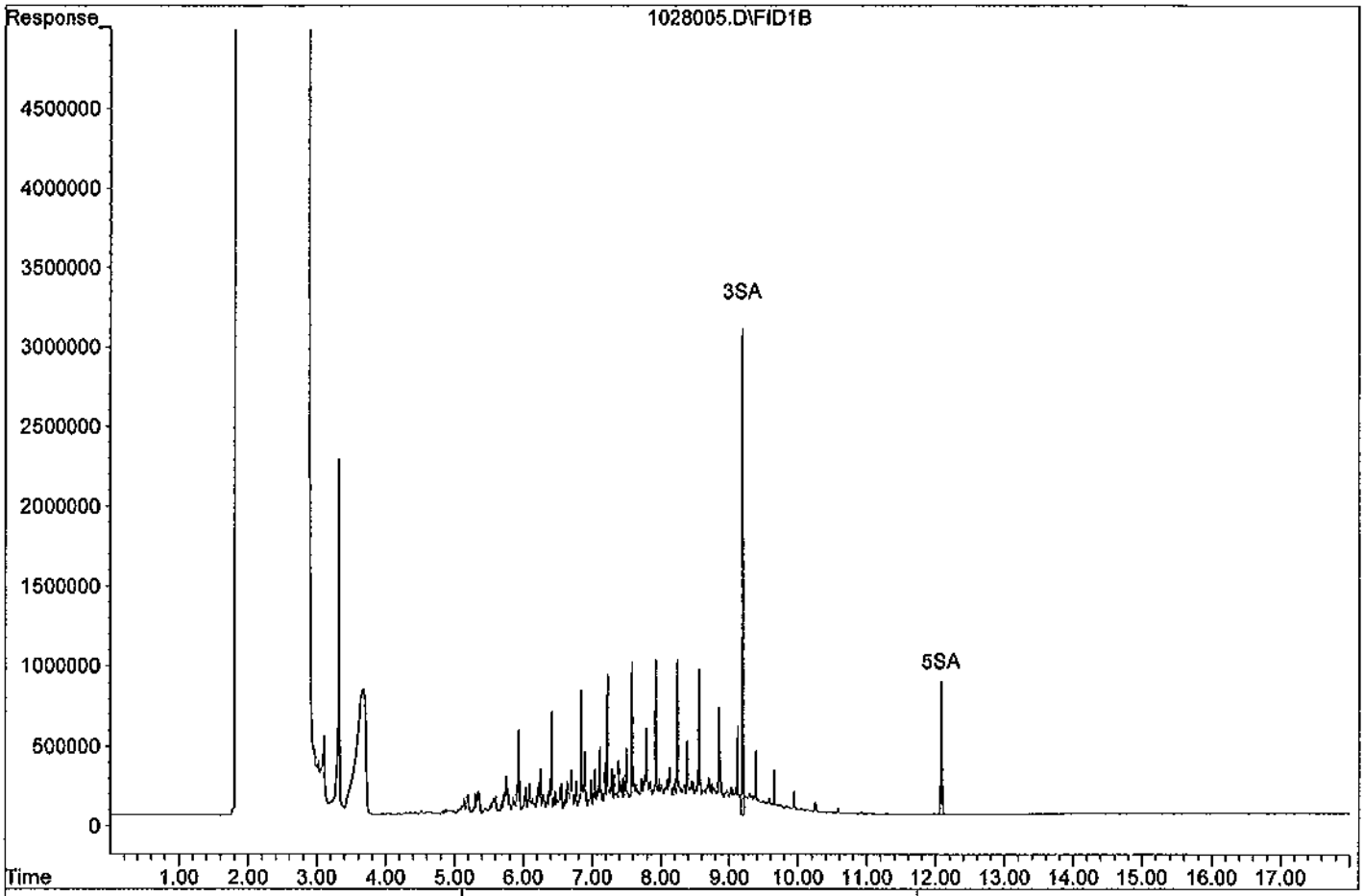
Method : G:\APOLLO\DATA\111028\TPH1028.M (Chemstation Integrator)
Title : Diesel
Last Update : Mon Oct 31 10:02:11 2011
Response via : Multiple Level Calibration

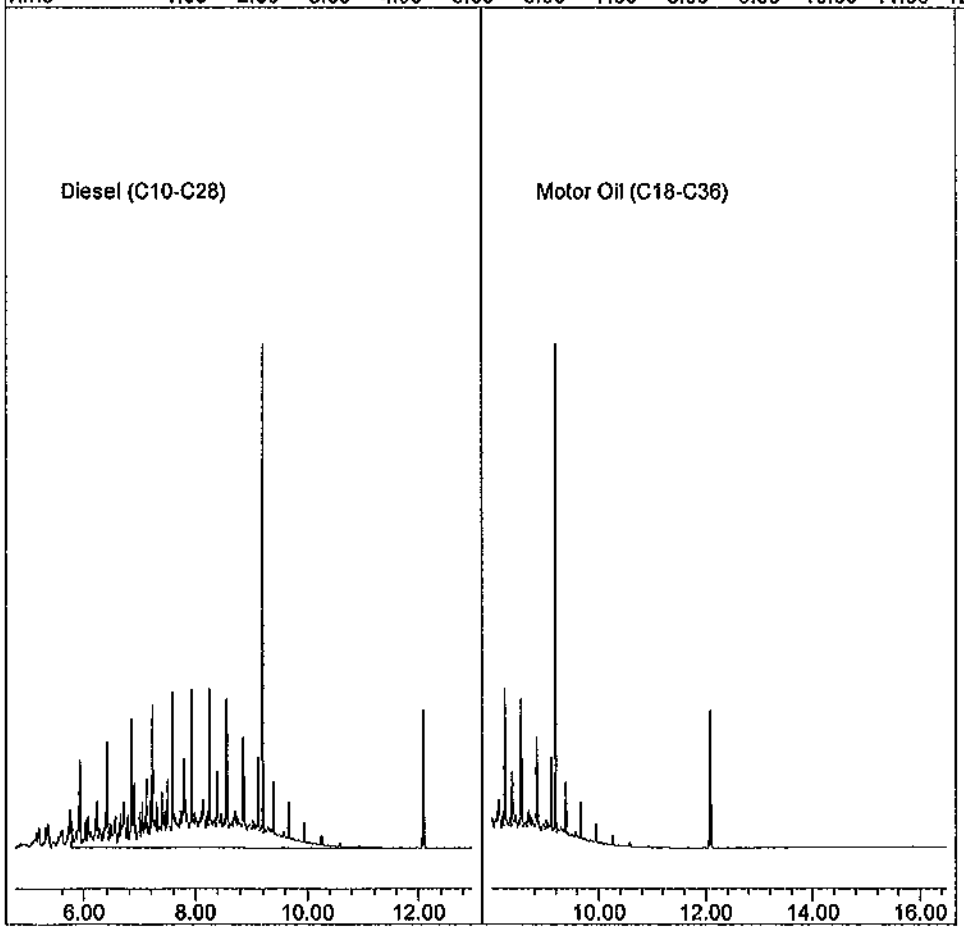
Volume Inj. : 2UL
Signal Phase : DB-5
Signal Info : FID02A

Compound	R.T.	Response	Conc Units
System Monitoring Compounds			
3) SA Not Used(S)	9.19	21988346	34.683 ppb
Surrogate Spike 30.000	Recovery =		115.61%
5) SA Not Used2(S)	12.09	10887525	43.308 ppb
Surrogate Spike 30.000	Recovery =		144.36%
Target Compounds			
1) HATM Diesel (C10-C28)	8.86	378076624	436.171 ppb

(f)=RT Delta > 1/2 Window 40 (m)=manual int.

Data File: G:\APOLLO\DATA\111028\1028005.D
Sample : DIESEL 400/1000





Quantitation Report (Not Reviewed)

Data File : G:\APOLLO\DATA\111028\1028006.D Vial: 6
Acq On : 10-28-11 10:59:35 Operator: LAC
Sample : DIESEL 600/1000 Inst : Apollo
Misc : Mix(A) Multiplr: 1.00
IntFile : events.e
Quant Time: Oct 28 13:43 2011 Quant Results File: TPH1028.RES

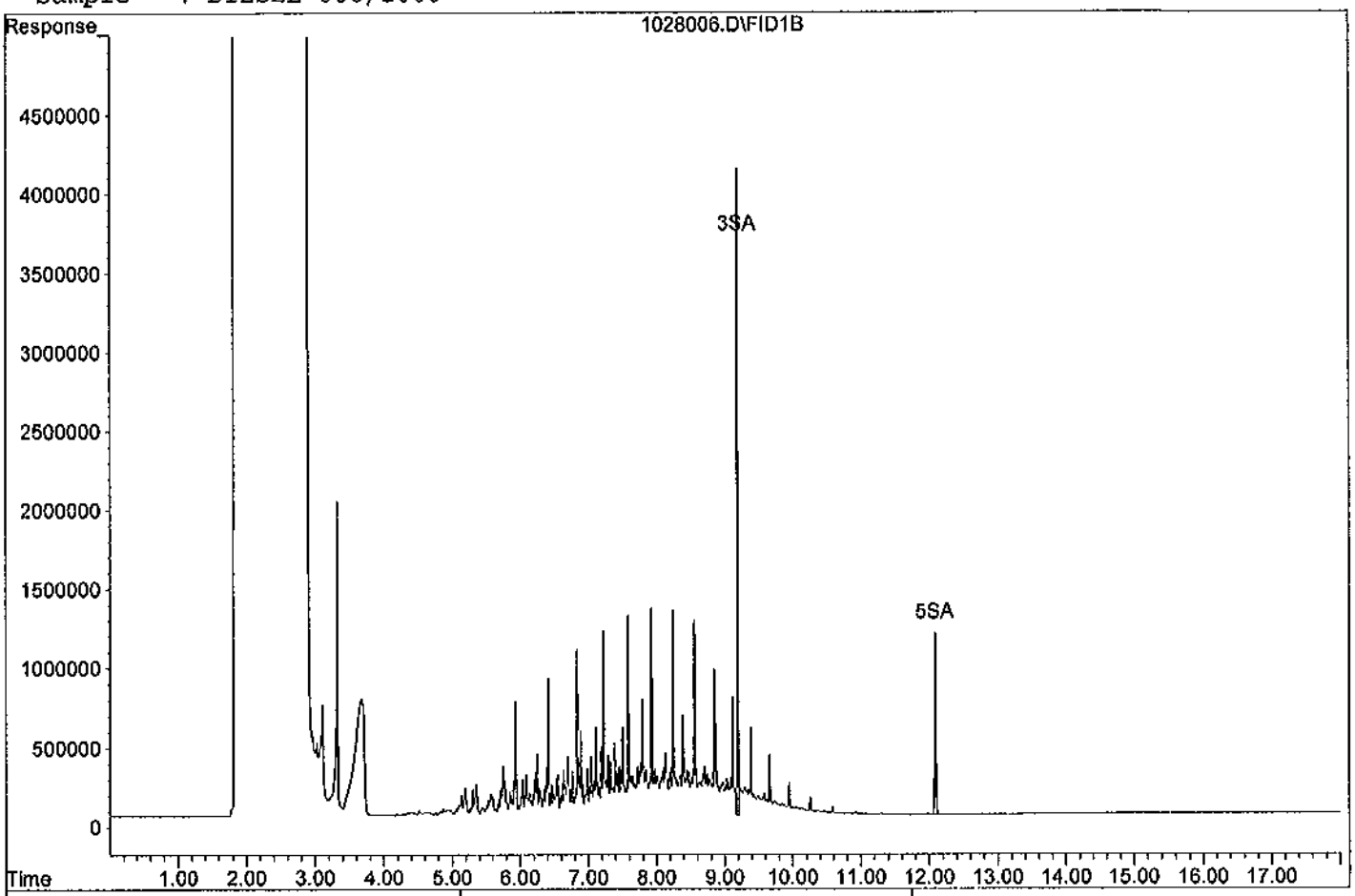
Method : G:\APOLLO\DATA\111028\TPH1028.M (Chemstation Integrator)
Title : Diesel
Last Update : Mon Oct 31 10:02:11 2011
Response via : Multiple Level Calibration

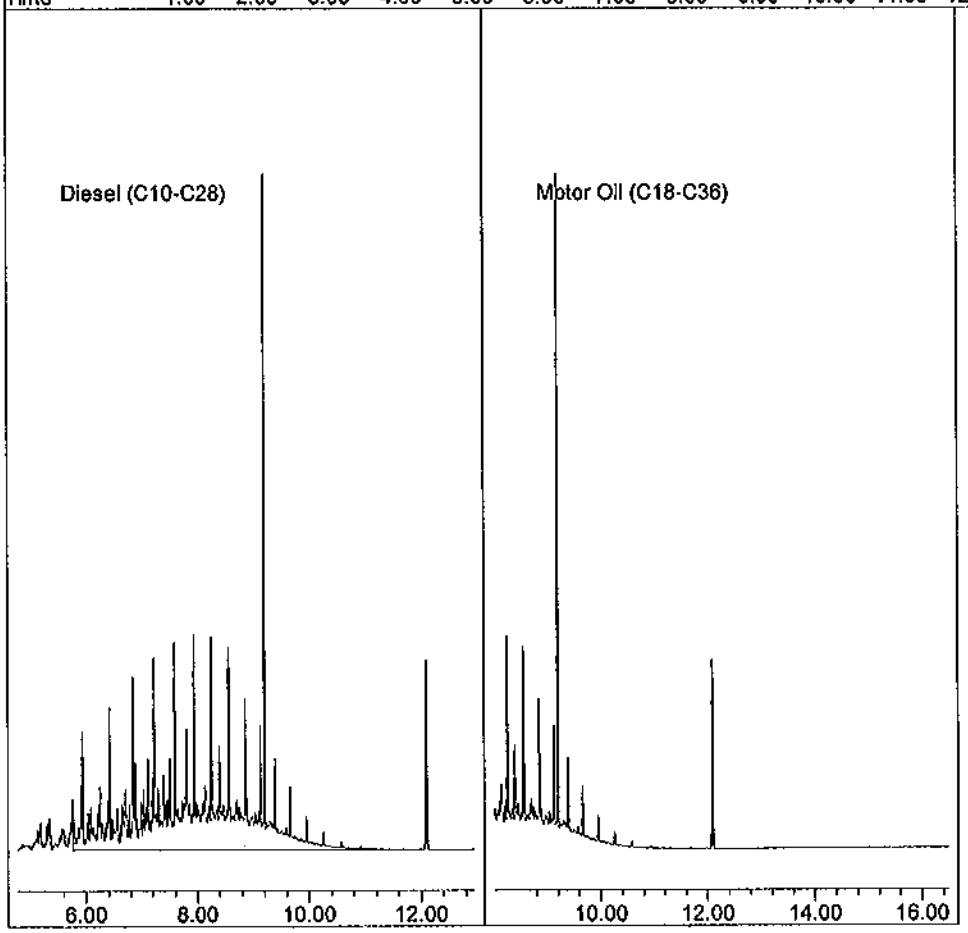
Volume Inj. : 2UL
Signal Phase : DB-5
Signal Info : FID02A

Compound	R.T.	Response	Conc Units
System Monitoring Compounds			
3) SA Not Used(S)	9.20	29970170	47.273 ppb
Surrogate Spike 30.000	Recovery =		157.58%
5) SA Not Used2(S)	12.09	15171855	60.906 ppb
Surrogate Spike 30.000	Recovery =		203.02%
Target Compounds			
1) HATM Diesel (C10-C28)	8.86	508771749	586.949 ppb

(f)=RT Delta > 1/2 Window 42 (m)=manual int.

Data File: G:\APOLLO\DATA\111028\1028006.D
Sample : DIESEL 600/1000





Quantitation Report (Not Reviewed)

Data File : G:\APOLLO\DATA\111028\1028007.D Vial: 7
Acq On : 10-28-11 11:23:49 Operator: LAC
Sample : DIESEL 800/1000 Inst : Apollo
Misc : Mix(A) Multiplr: 1.00
IntFile : events.e
Quant Time: Oct 28 13:43 2011 Quant Results File: TPH1028.RES

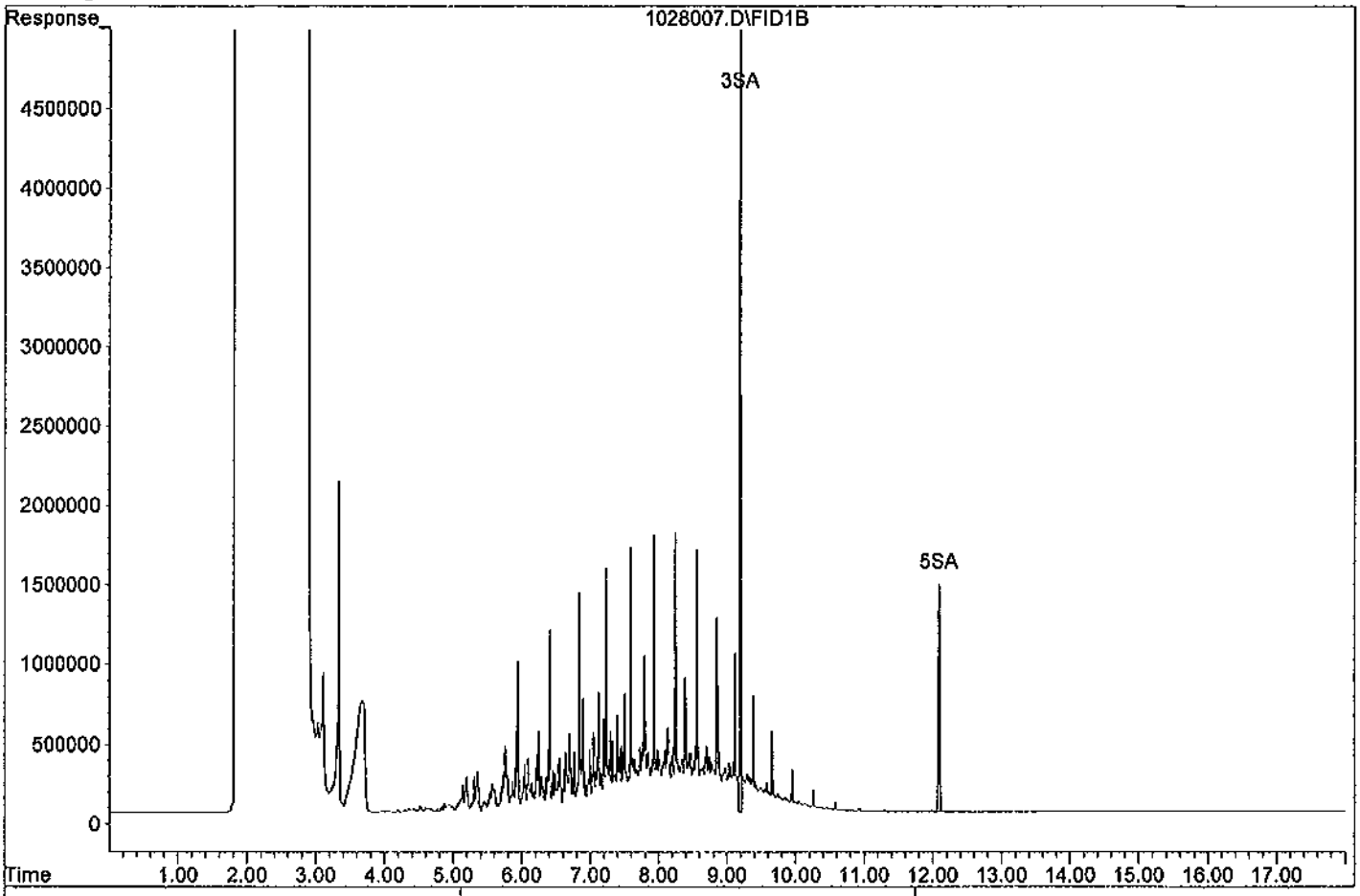
Method : G:\APOLLO\DATA\111028\TPH1028.M (Chemstation Integrator)
Title : Diesel
Last Update : Mon Oct 31 10:02:11 2011
Response via : Multiple Level Calibration

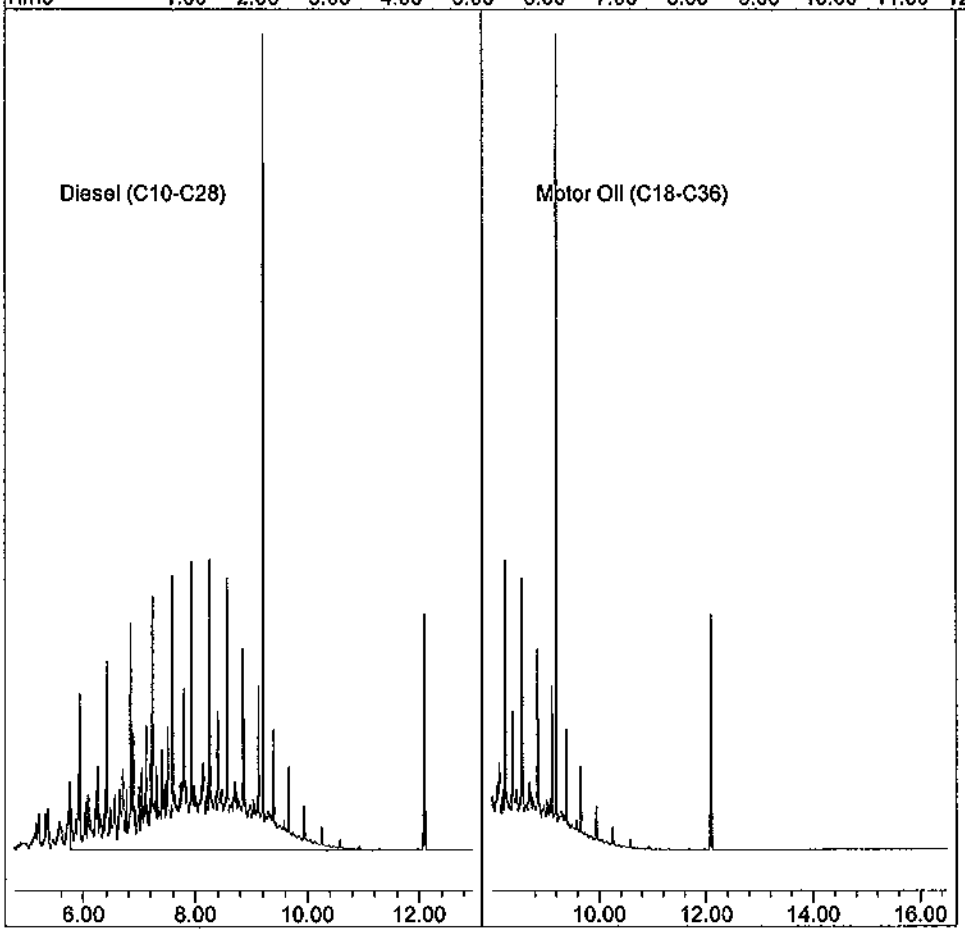
Volume Inj. : 2UL
Signal Phase : DB-5
Signal Info : FID02A

Compound	R.T.	Response	Conc Units
System Monitoring Compounds			
3) SA Not Used(S)	9.20	40429014	63.770 ppb
Surrogate Spike 30.000	Recovery =		212.57%
5) SA Not Used2(S)	12.10	19720236	79.588 ppb
Surrogate Spike 30.000	Recovery =		265.29%
Target Compounds			
1) HATM Diesel (C10-C28)	8.86	680806039	785.417 ppb

(f)=RT Delta > 1/2 Window 44 (m)=manual int.

Data File: G:\APOLLO\DATA\111028\1028007.D
Sample : DIESEL 800/1000





Quantitation Report (Not Reviewed)

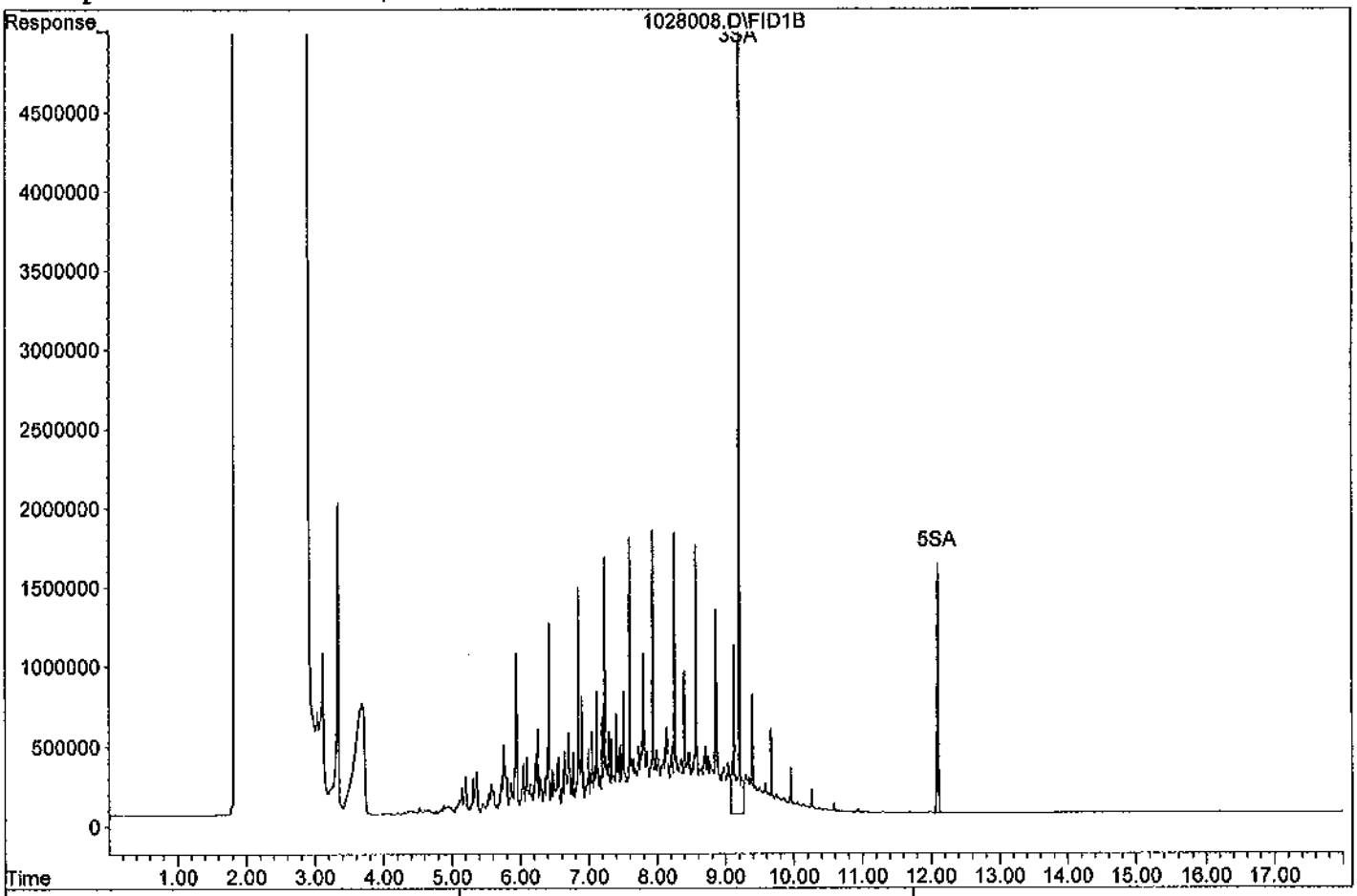
Data File : G:\APOLLO\DATA\111028\1028008.D Vial: 8
Acq On : 10-28-11 11:48:05 Operator: LAC
Sample : DIESEL 1000/1000 Inst : Apollo
Misc : Mix(A) Multiplr: 1.00
IntFile : events.e
Quant Time: Oct 28 13:43 2011 Quant Results File: TPH1028.RES

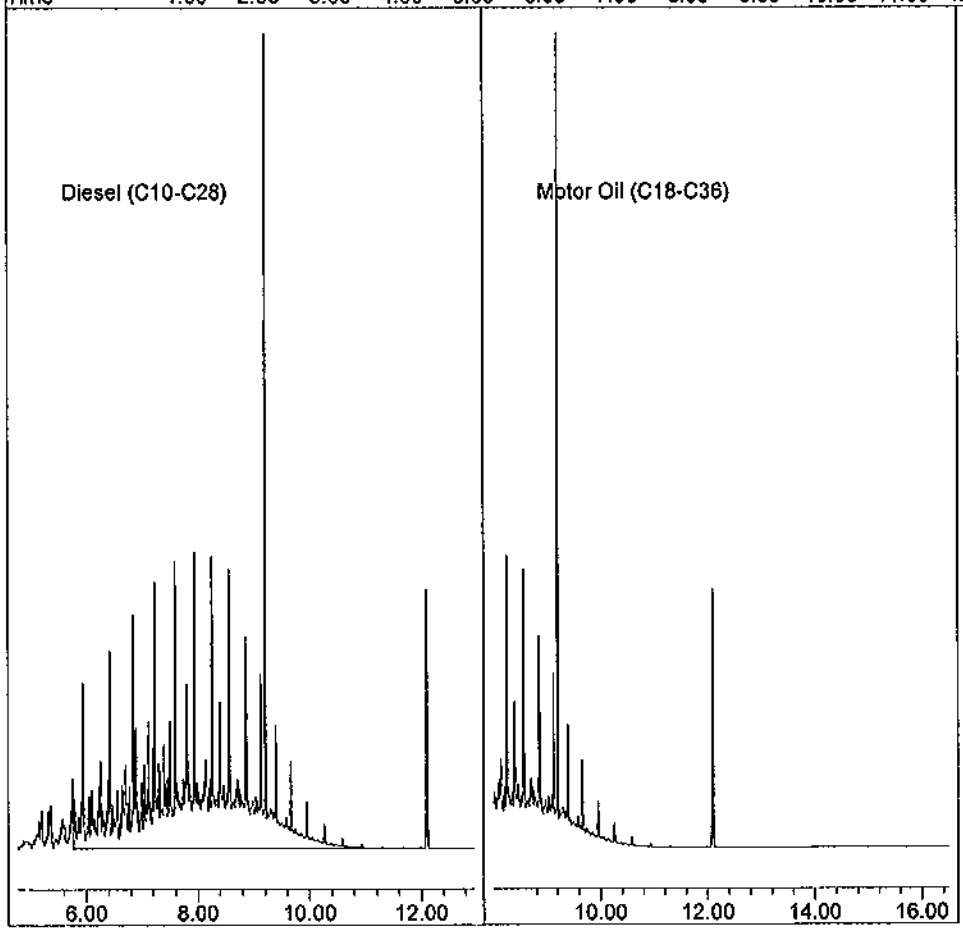
Method : G:\APOLLO\DATA\111028\TPH1028.M (Chemstation Integrator)
Title : Diesel
Last Update : Mon Oct 31 10:02:11 2011
Response via : Multiple Level Calibration

Volume Inj. : 2UL
Signal Phase : DB-5
Signal Info : FID02A

Compound	R.T.	Response	Conc Units
System Monitoring Compounds			
3) SA Not Used(S)	9.20	68048921	107.335 ppb
Surrogate Spike 30.000	Recovery =		357.78%
5) SA Not Used2(S)	12.10	21357059	86.311 ppb
Surrogate Spike 30.000	Recovery =		287.70%
Target Compounds			
1) HATM Diesel (C10-C28)	8.86	680836698	785.453 ppb

Data File: G:\APOLLO\DATA\111028\1028008.D
Sample : DIESEL 1000/1000





Quantitation Report (Not Reviewed)

Data File : G:\APOLLO\DATA\111028\1028009.D Vial: 9
Acq On : 10-28-11 12:12:27 Operator: LAC
Sample : MOTOR OIL 50/1000 10/28/11 Inst : Apollo
Misc : Mix(B) Multiplr: 1.00
IntFile : events.e
Quant Time: Oct 28 13:43 2011 Quant Results File: TPH1028.RES

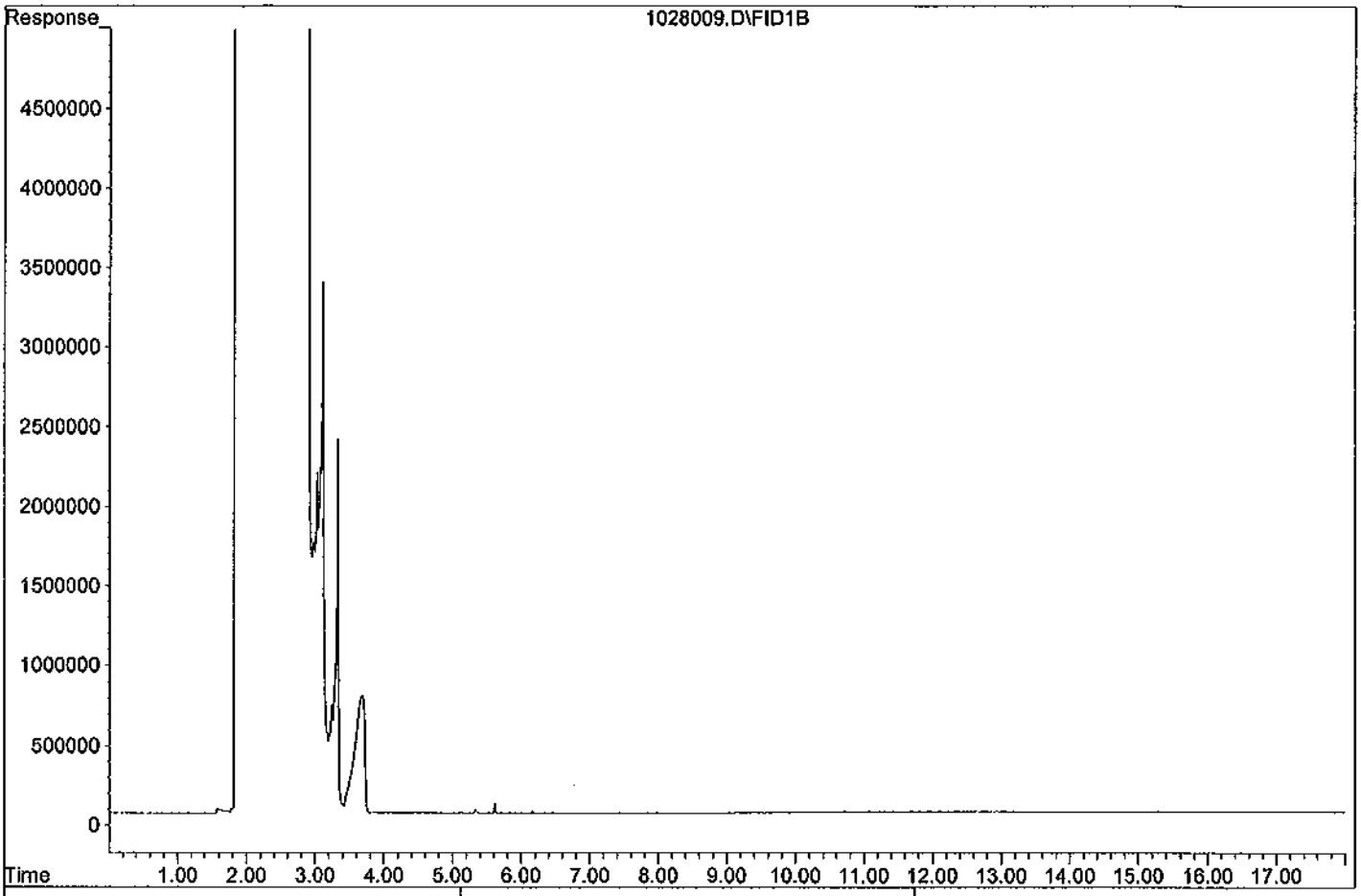
Method : G:\APOLLO\DATA\111028\TPH1028.M (Chemstation Integrator)
Title : Diesel
Last Update : Mon Oct 31 10:02:11 2011
Response via : Multiple Level Calibration

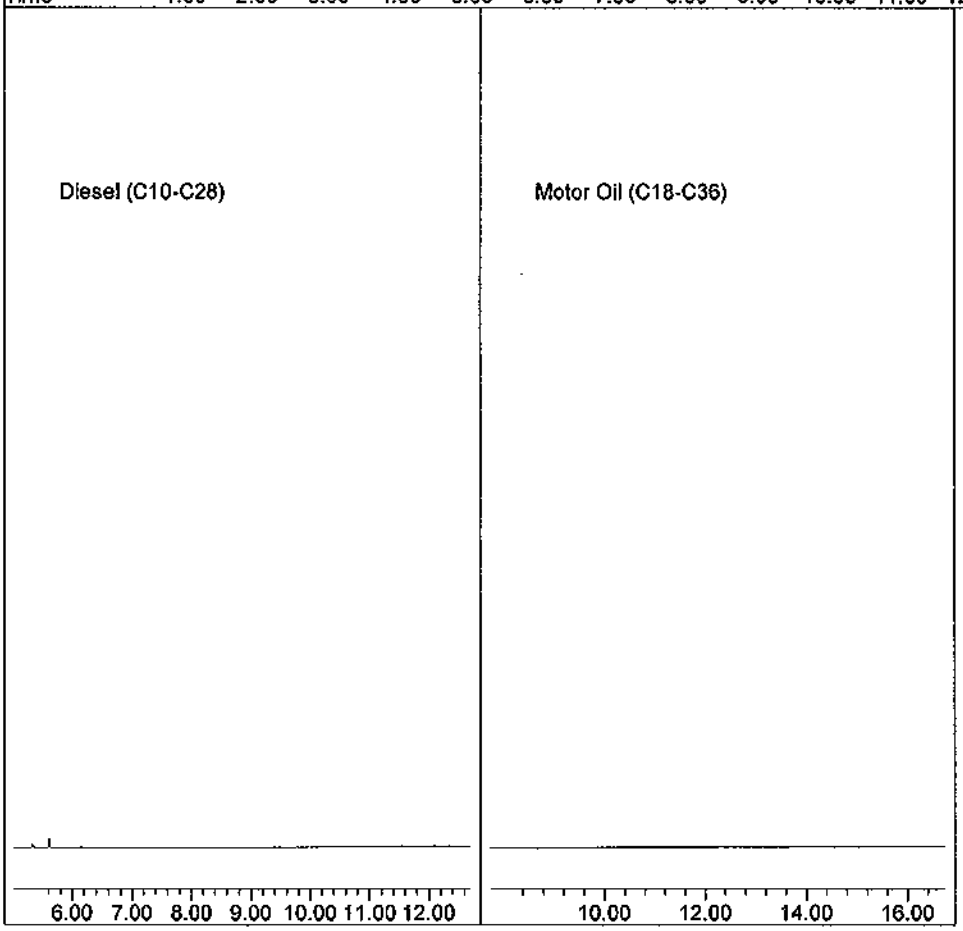
Volume Inj. : 2UL
Signal Phase : DB-5
Signal Info : FID02A

Compound	R.T.	Response	Conc Units
System Monitoring Compounds			
Target Compounds			
2) HBTM Motor Oil (C18-C36)	12.25	17842259	49.322 ppb

(f)=RT Delta > 1/2 Window 48 (m)=manual int.

Data File: G:\APOLLO\DATA\111028\1028009.D
Sample : MOTOR OIL 50/1000 10/28/11





Quantitation Report (Not Reviewed)

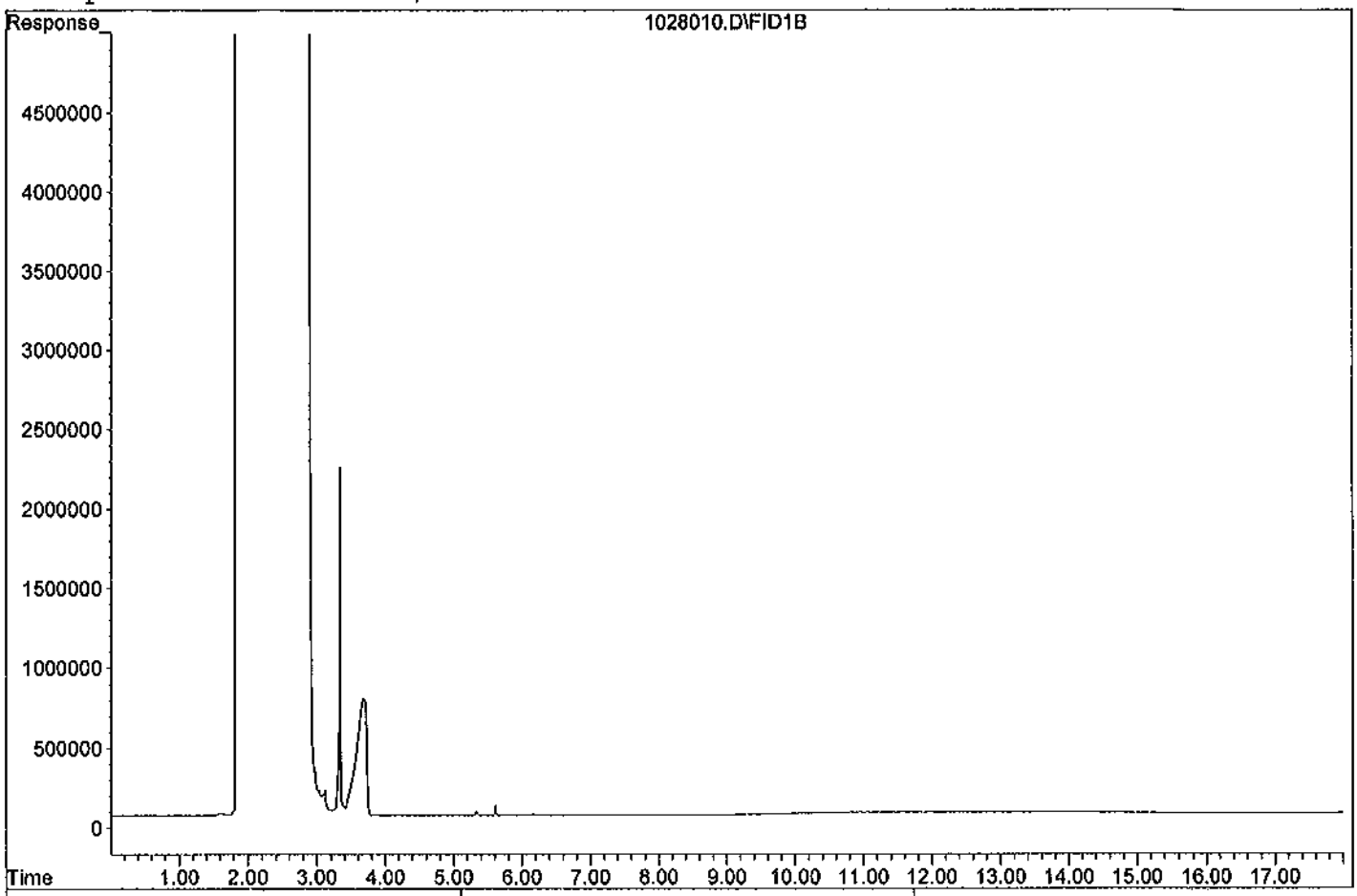
Data File : G:\APOLLO\DATA\111028\1028010.D Vial: 10
Acq On : 10-28-11 12:36:20 Operator: LAC
Sample : MOTOR OIL 100/1000 Inst : Apollo
Misc : Mix(B) Multiplr: 1.00
IntFile : events.e
Quant Time: Oct 28 13:43 2011 Quant Results File: TPH1028.RES

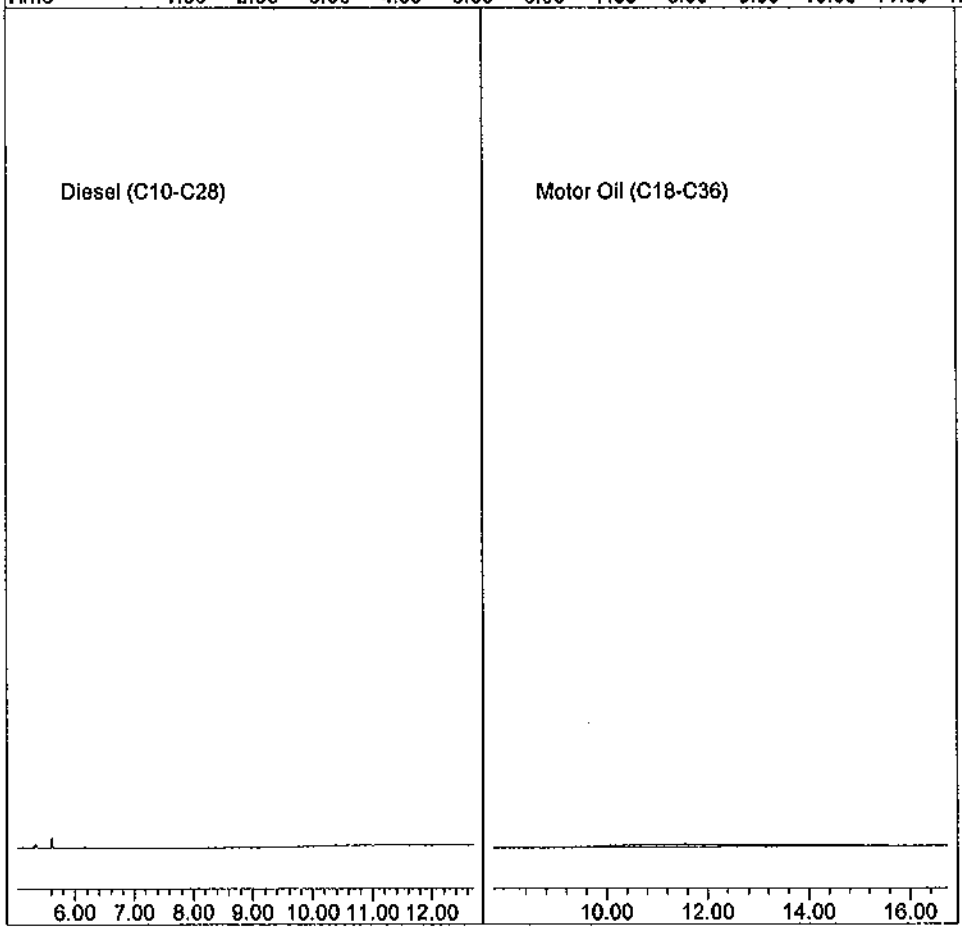
Method : G:\APOLLO\DATA\111028\TPH1028.M (Chemstation Integrator)
Title : Diesel
Last Update : Mon Oct 31 10:02:11 2011
Response via : Multiple Level Calibration

Volume Inj. : 2UL
Signal Phase : DB-5
Signal Info : FID02A

Compound	R.T.	Response	Conc Units
System Monitoring Compounds			
Target Compounds			
2) HBTM Motor Oil (C18-C36)	12.25	36743279	101.570 ppb

Data File: G:\APOLLO\DATA\111028\1028010.D
Sample : MOTOR OIL 100/1000





Quantitation Report (Not Reviewed)

Data File : G:\APOLLO\DATA\111028\1028011.D Vial: 11
Acq On : 10-28-11 13:00:16 Operator: LAC
Sample : MOTOR OIL 400/1000 Inst : Apollo
Misc : Mix(B) Multiplr: 1.00
IntFile : events.e
Quant Time: Oct 28 13:43 2011 Quant Results File: TPH1028.RES

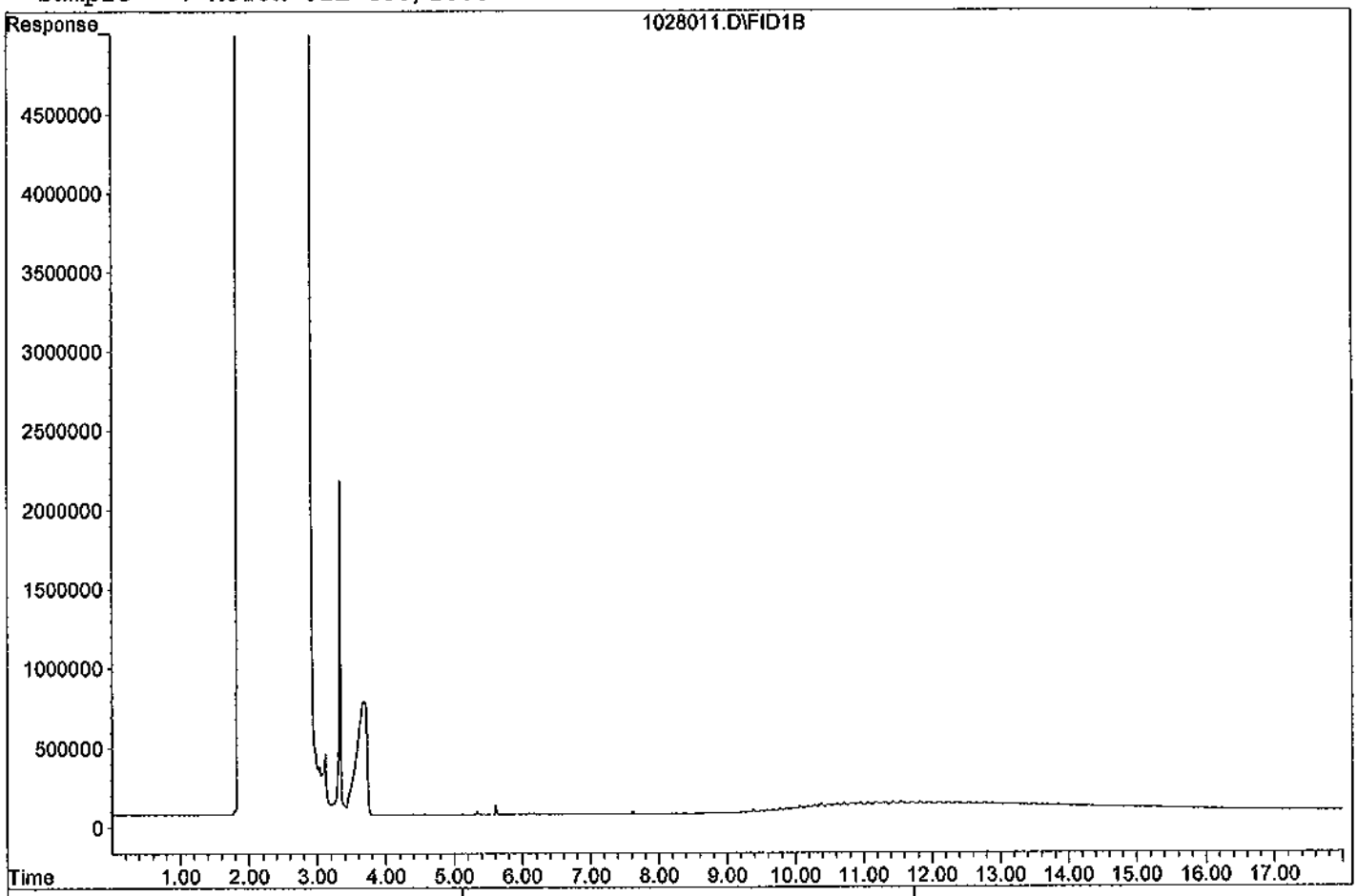
Method : G:\APOLLO\DATA\111028\TPH1028.M (Chemstation Integrator)
Title : Diesel
Last Update : Mon Oct 31 10:02:11 2011
Response via : Multiple Level Calibration

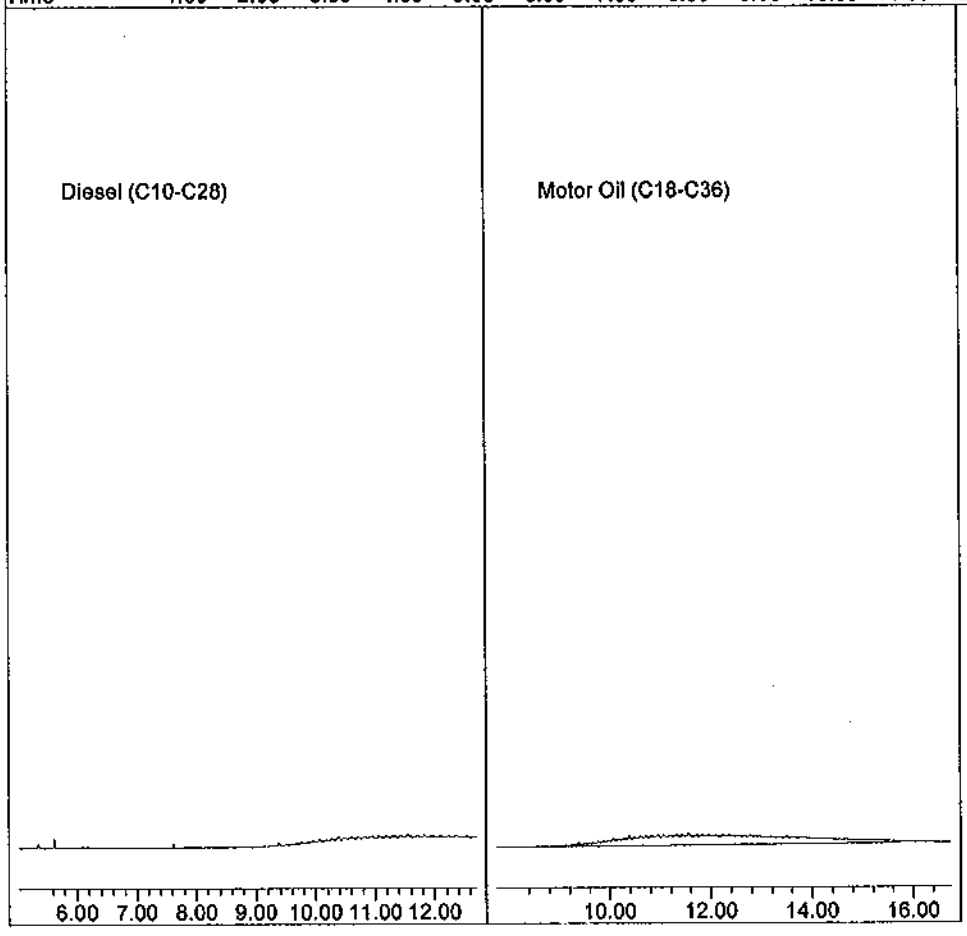
Volume Inj. : 2UL
Signal Phase : DB-5
Signal Info : FID02A

Compound	R.T.	Response	Conc Units
System Monitoring Compounds			
Target Compounds			
2) HBTM Motor Oil (C18-C36)	12.25	147050915	406.495 ppb

(f)=RT Delta > 1/2 Window 52 (m)=manual int.

Data File: G:\APOLLO\DATA\111028\1028011.D
Sample : MOTOR OIL 400/1000





Quantitation Report (Not Reviewed)

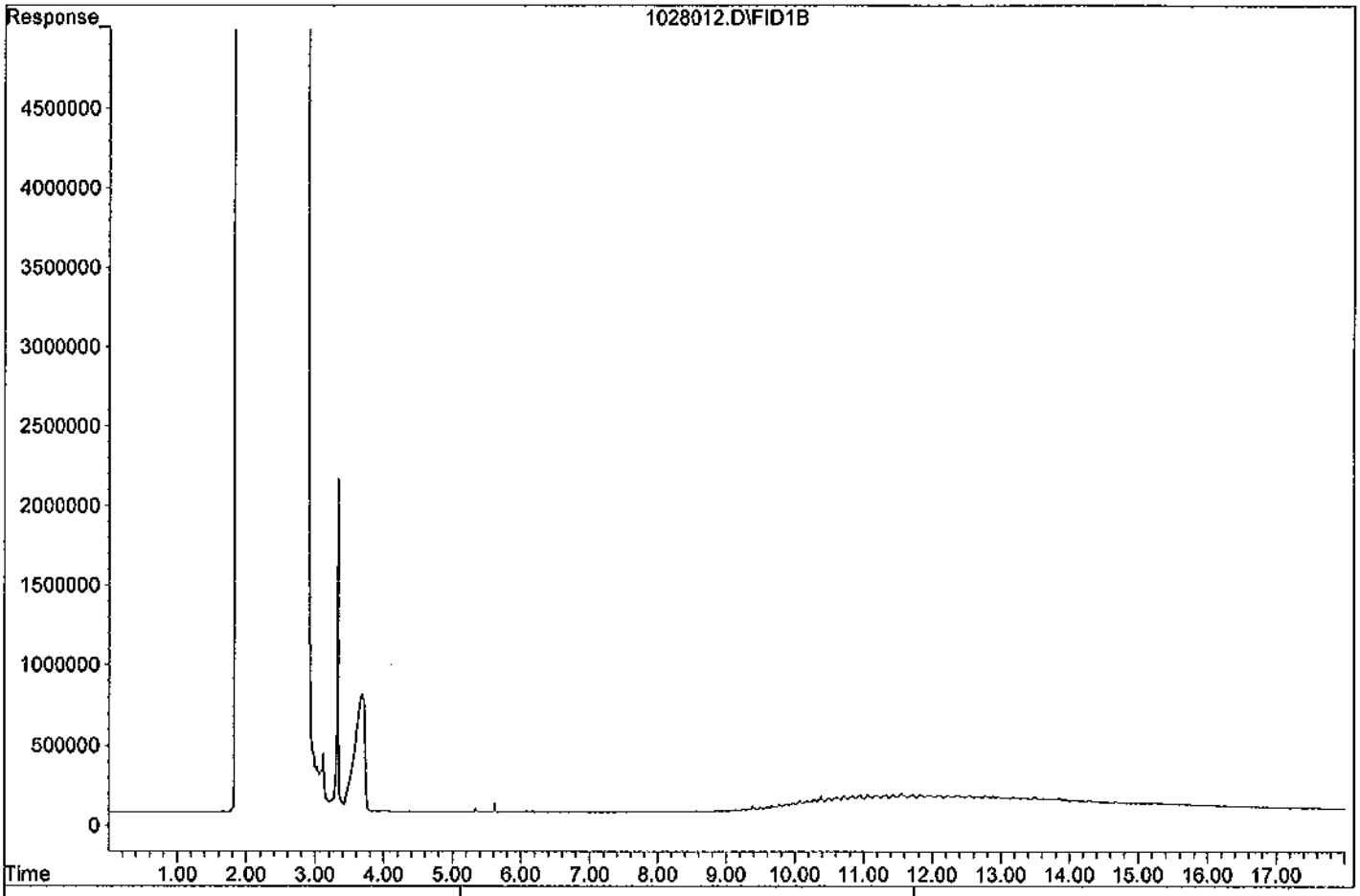
Data File : G:\APOLLO\DATA\111028\1028012.D Vial: 12
Acq On : 10-28-11 13:24:39 Operator: LAC
Sample : MOTOR OIL 600/1000 Inst : Apollo
Misc : Mix(B) Multiplr: 1.00
IntFile : events.e
Quant Time: Oct 28 13:43 2011 Quant Results File: TPH1028.RES

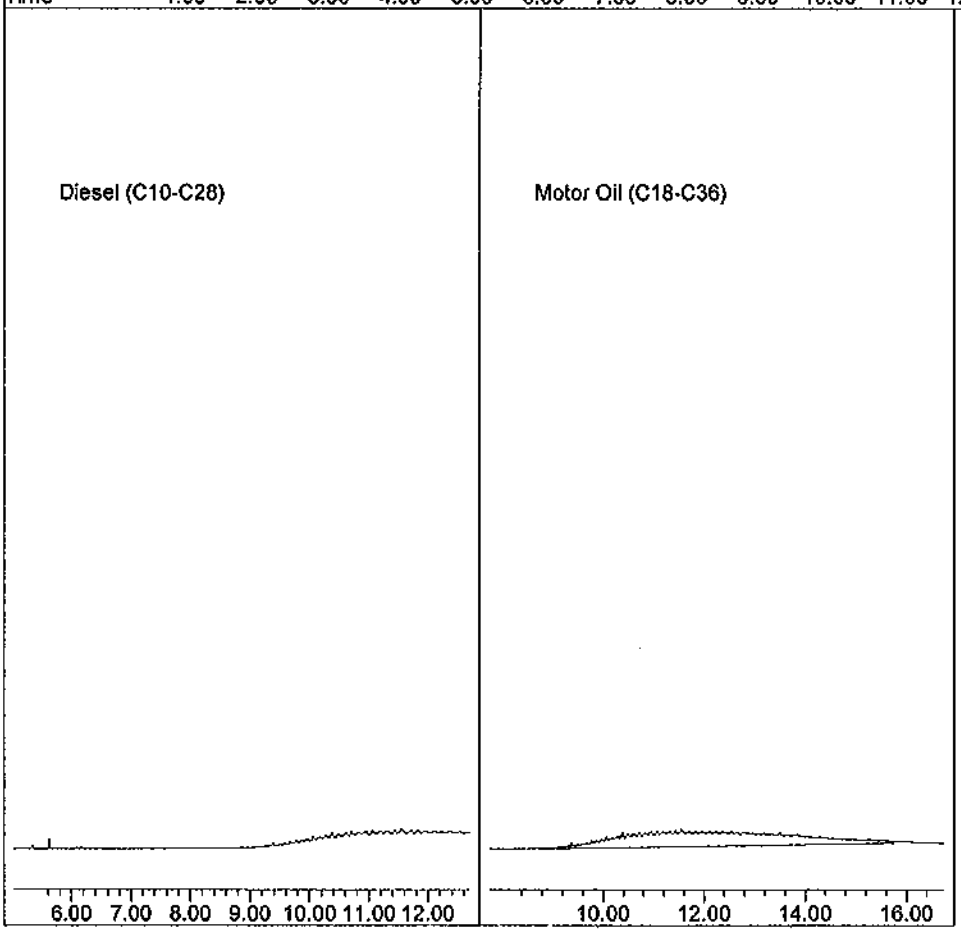
Method : G:\APOLLO\DATA\111028\TPH1028.M (Chemstation Integrator)
Title : Diesel
Last Update : Mon Oct 31 10:02:11 2011
Response via : Multiple Level Calibration

Volume Inj. : 2UL
Signal Phase : DB-5
Signal Info : FID02A

Compound	R.T.	Response	Conc Units
System Monitoring Compounds			
Target Compounds			
2) HBTM Motor Oil (C18-C36)	12.25	216778154	599.242 ppb

Data File: G:\APOLLO\DATA\111028\1028012.D
Sample : MOTOR OIL 600/1000





Quantitation Report (Not Reviewed)

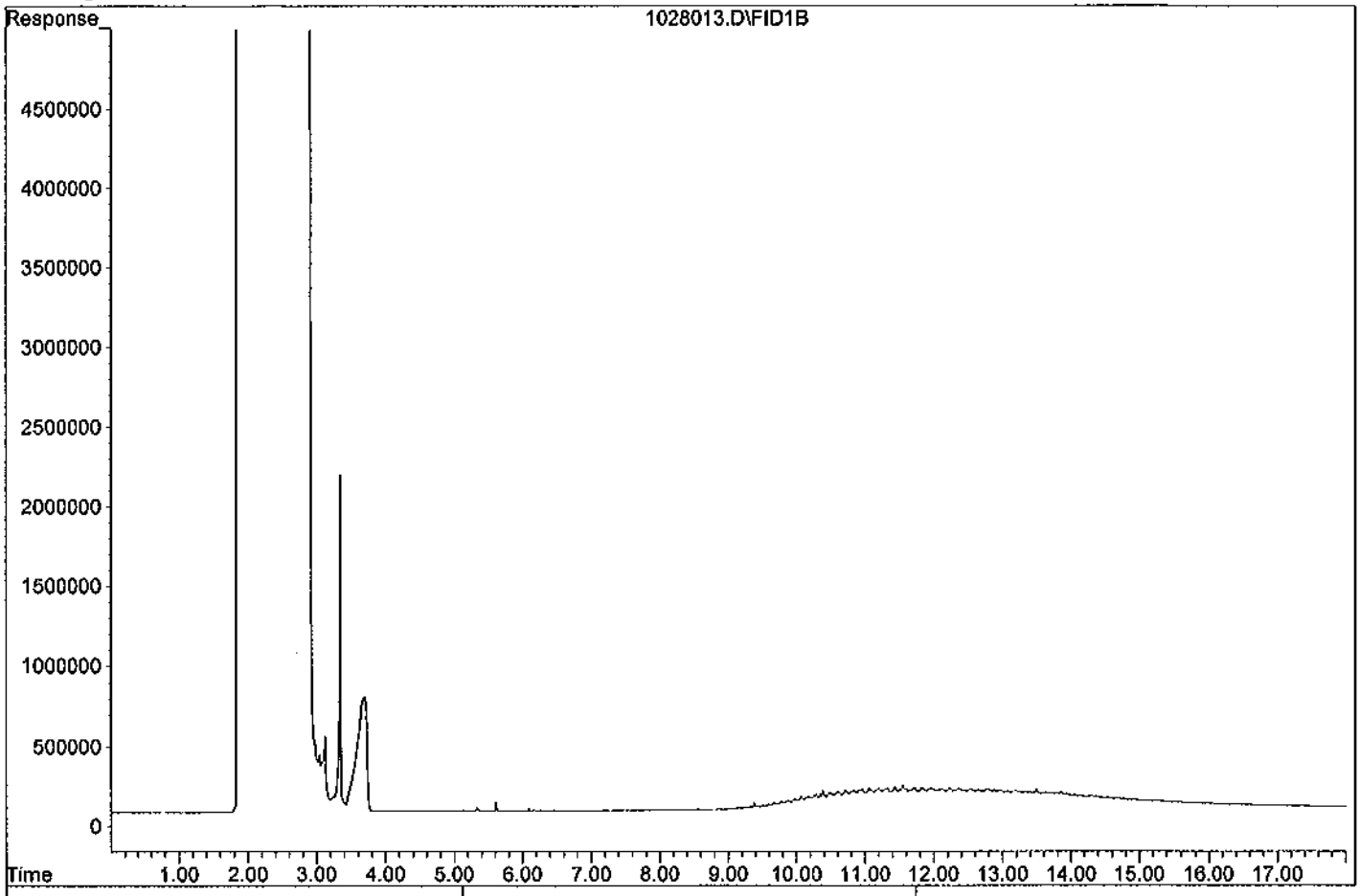
Data File : G:\APOLLO\DATA\111028\1028013.D Vial: 13
Acq On : 10-28-11 13:48:43 Operator: LAC
Sample : MOTOR OIL 800/1000 Inst : Apollo
Misc : Mix(B) Multiplr: 1.00
IntFile : events.e
Quant Time: Oct 28 13:43 2011 Quant Results File: TPH1028.RES

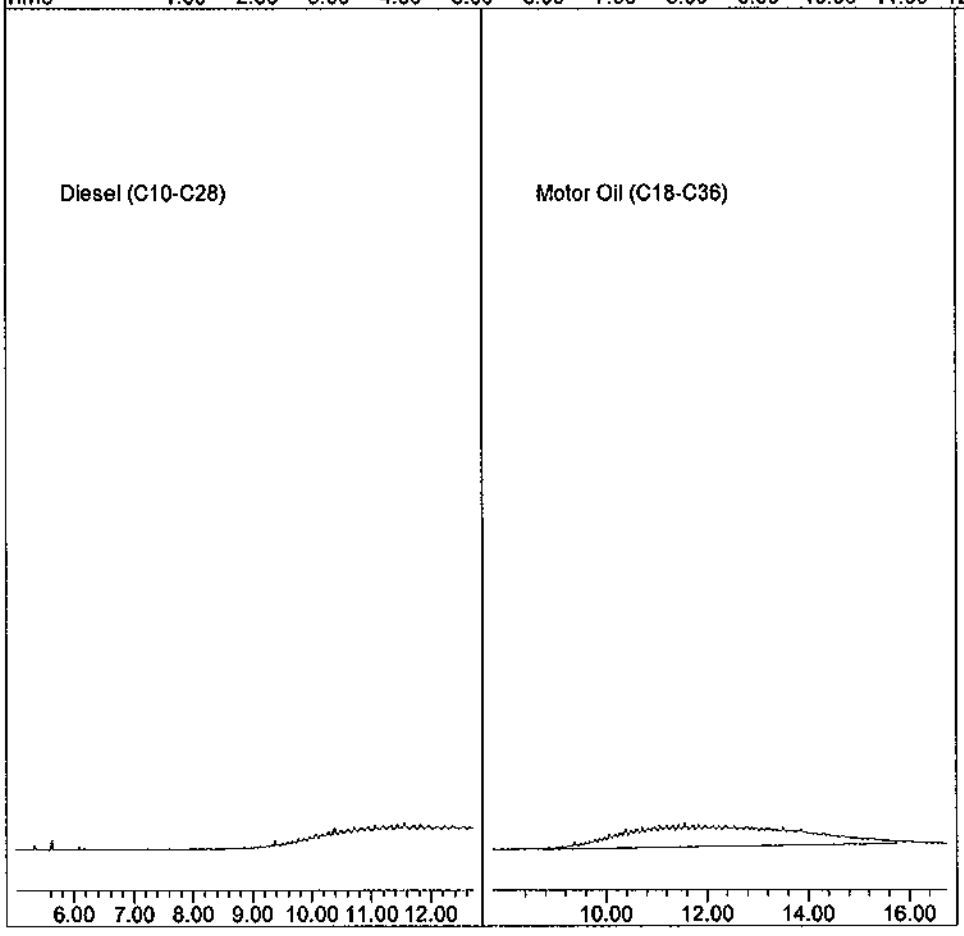
Method : G:\APOLLO\DATA\111028\TPH1028.M (Chemstation Integrator)
Title : Diesel
Last Update : Mon Oct 31 10:02:11 2011
Response via : Multiple Level Calibration

Volume Inj. : 2UL
Signal Phase : DB-5
Signal Info : FID02A

Compound	R.T.	Response	Conc Units
System Monitoring Compounds			
Target Compounds			
2) HBTM Motor Oil (C18-C36)	12.25	303785051	839.757 ppb

Data File: G:\APOLLO\DATA\111028\1028013.D
Sample : MOTOR OIL 800/1000





Quantitation Report (Not Reviewed)

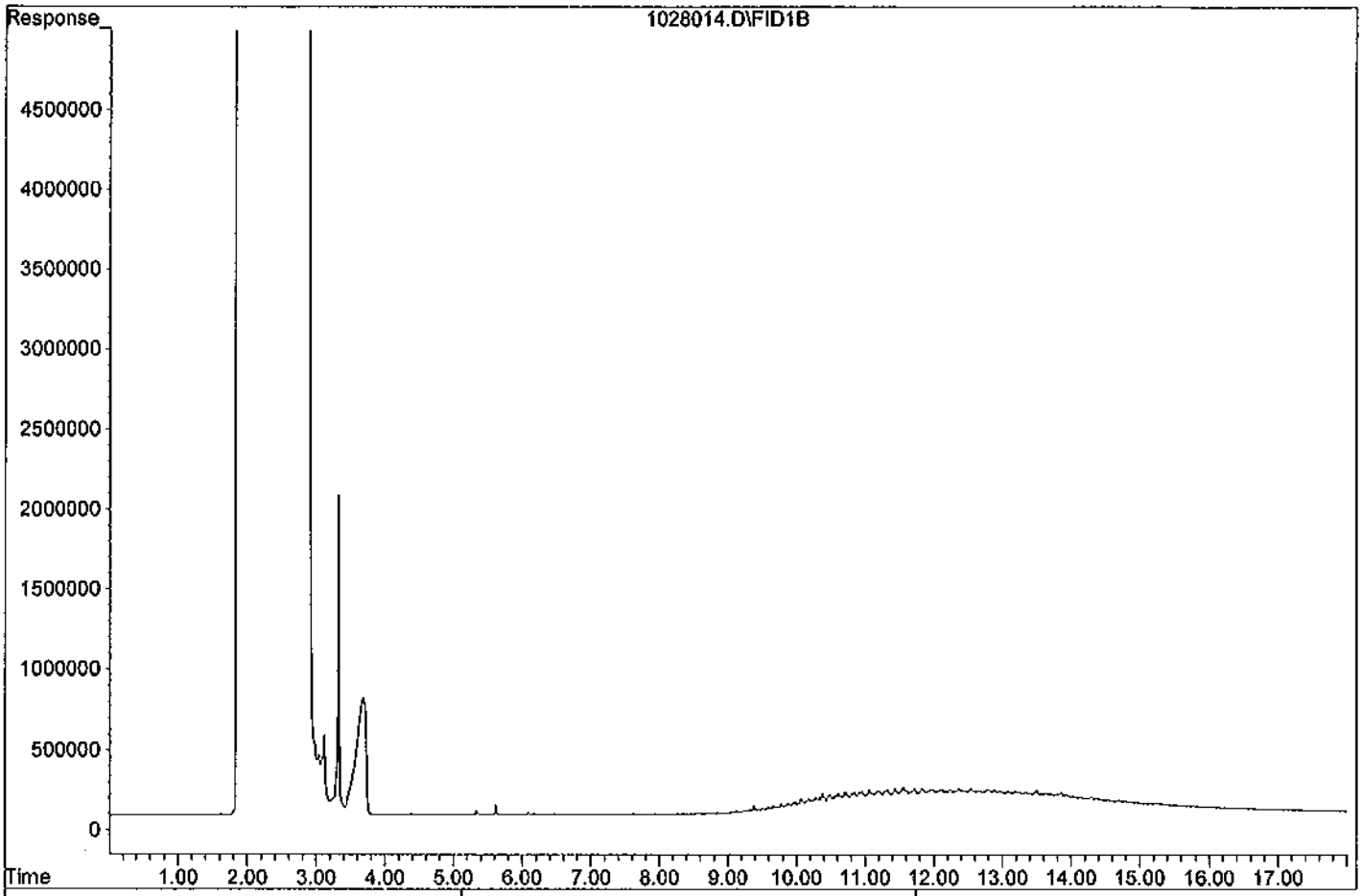
Data File : G:\APOLLO\DATA\111028\1028014.D Vial: 14
Acq On : 10-28-11 14:13:14 Operator: LAC
Sample : MOTOR OIL 1000/1000 Inst : Apollo
Misc : Mix(B) Multiplr: 1.00
IntFile : events.e
Quant Time: Oct 28 13:43 2011 Quant Results File: TPH1028.RES

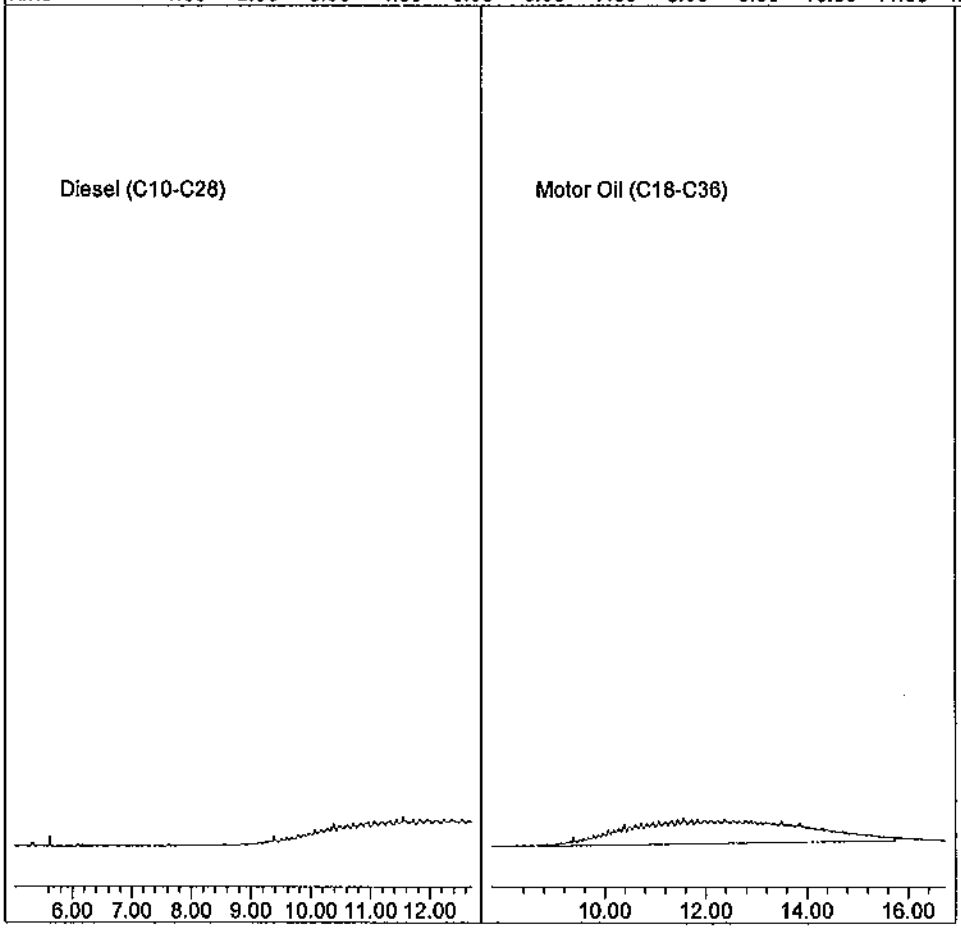
Method : G:\APOLLO\DATA\111028\TPH1028.M (Chemstation Integrator)
Title : Diesel
Last Update : Mon Oct 31 10:02:11 2011
Response via : Multiple Level Calibration

Volume Inj. : 2UL
Signal Phase : DB-5
Signal Info : FID02A

Compound	R.T.	Response	Conc Units
System Monitoring Compounds			
Target Compounds			
2) HBTM Motor Oil (C18-C36)	12.25	342332944	946.315 ppb

Data File: G:\APOLLO\DATA\111028\1028014.D
Sample : MOTOR OIL 1000/1000





Quantitation Report (Not Reviewed)

Data File : G:\APOLLO\DATA\111028\1028016.D Vial: 16
Acq On : 10-28-11 15:01:44 Operator: LAC
Sample : THC SURR 10/1000 10/28/11 Inst : Apollo
Misc : Mix(C) Multiplr: 1.00
IntFile : events.e
Quant Time: Oct 31 9:01 2011 Quant Results File: TPH1028.RES

Method : G:\APOLLO\DATA\111028\TPH1028.M (Chemstation Integrator)
Title : Diesel
Last Update : Mon Oct 31 10:02:11 2011
Response via : Multiple Level Calibration

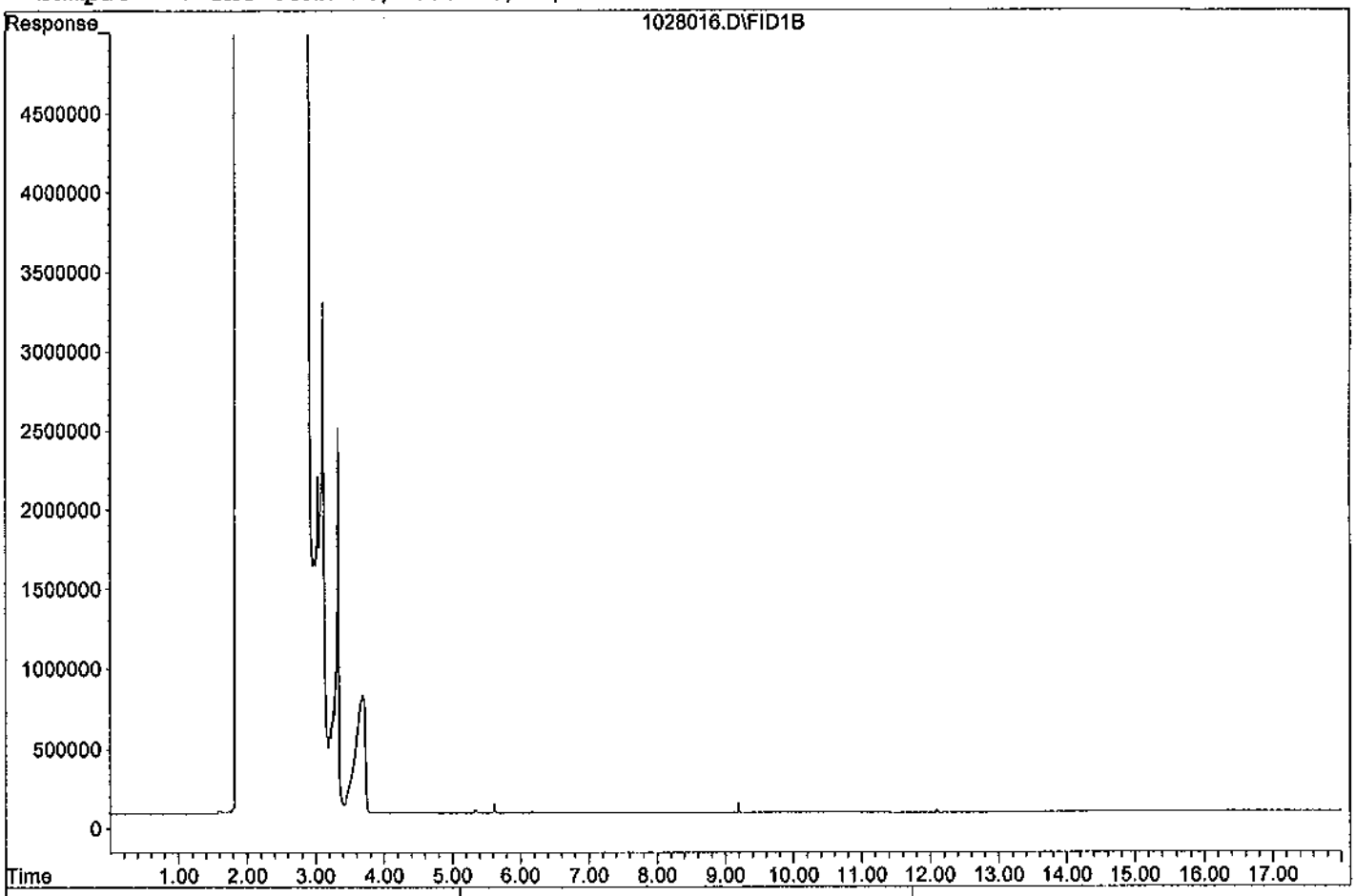
Volume Inj. : 2UL
Signal Phase : DB-5
Signal Info : FID02A

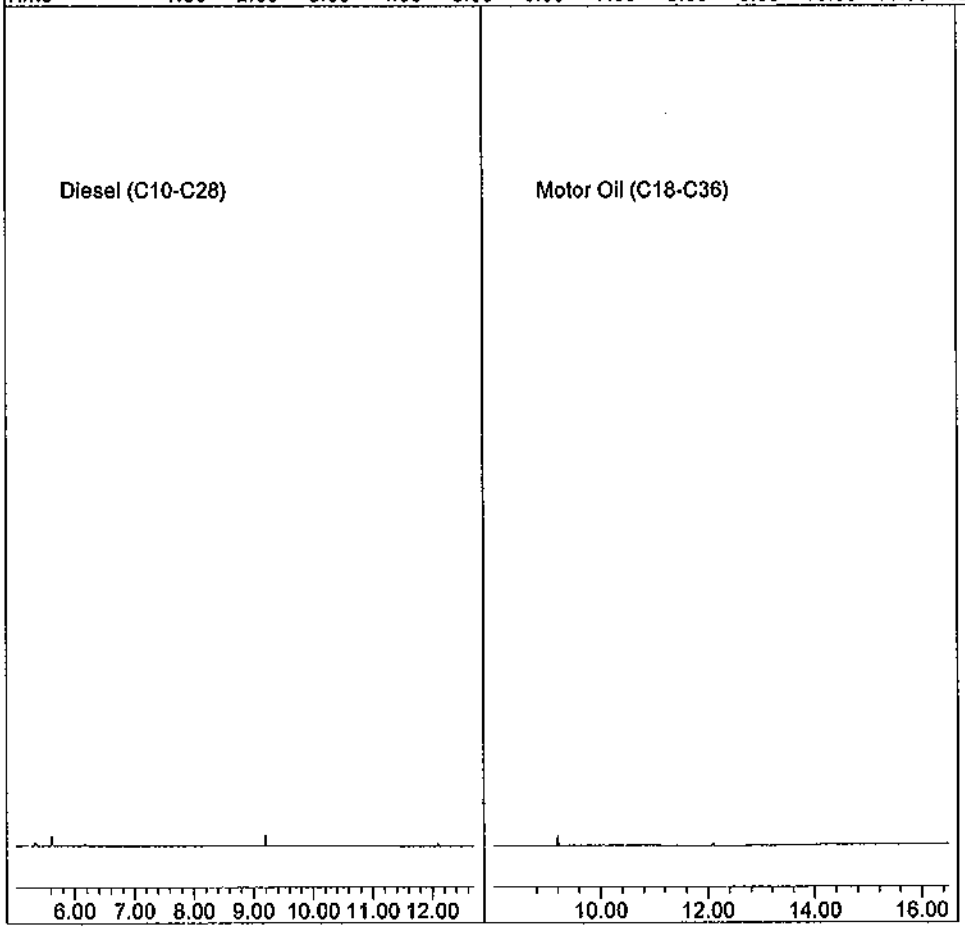
Compound	R.T.	Response	Conc Units

System Monitoring Compounds

Target Compounds

Data File: G:\APOLLO\DATA\111028\1028016.D
Sample : THC SURR 10/1000 10/28/11





Quantitation Report (Not Reviewed)

Data File : G:\APOLLO\DATA\111028\1028017.D Vial: 17
Acq On : 10-28-11 15:25:58 Operator: LAC
Sample : THC SURR 100/1000 Inst : Apollo
Misc : Mix(C) Multiplr: 1.00
IntFile : events.e
Quant Time: Oct 31 9:01 2011 Quant Results File: TPH1028.RES

Method : G:\APOLLO\DATA\111028\TPH1028.M (Chemstation Integrator)
Title : Diesel
Last Update : Mon Oct 31 10:02:11 2011
Response via : Multiple Level Calibration

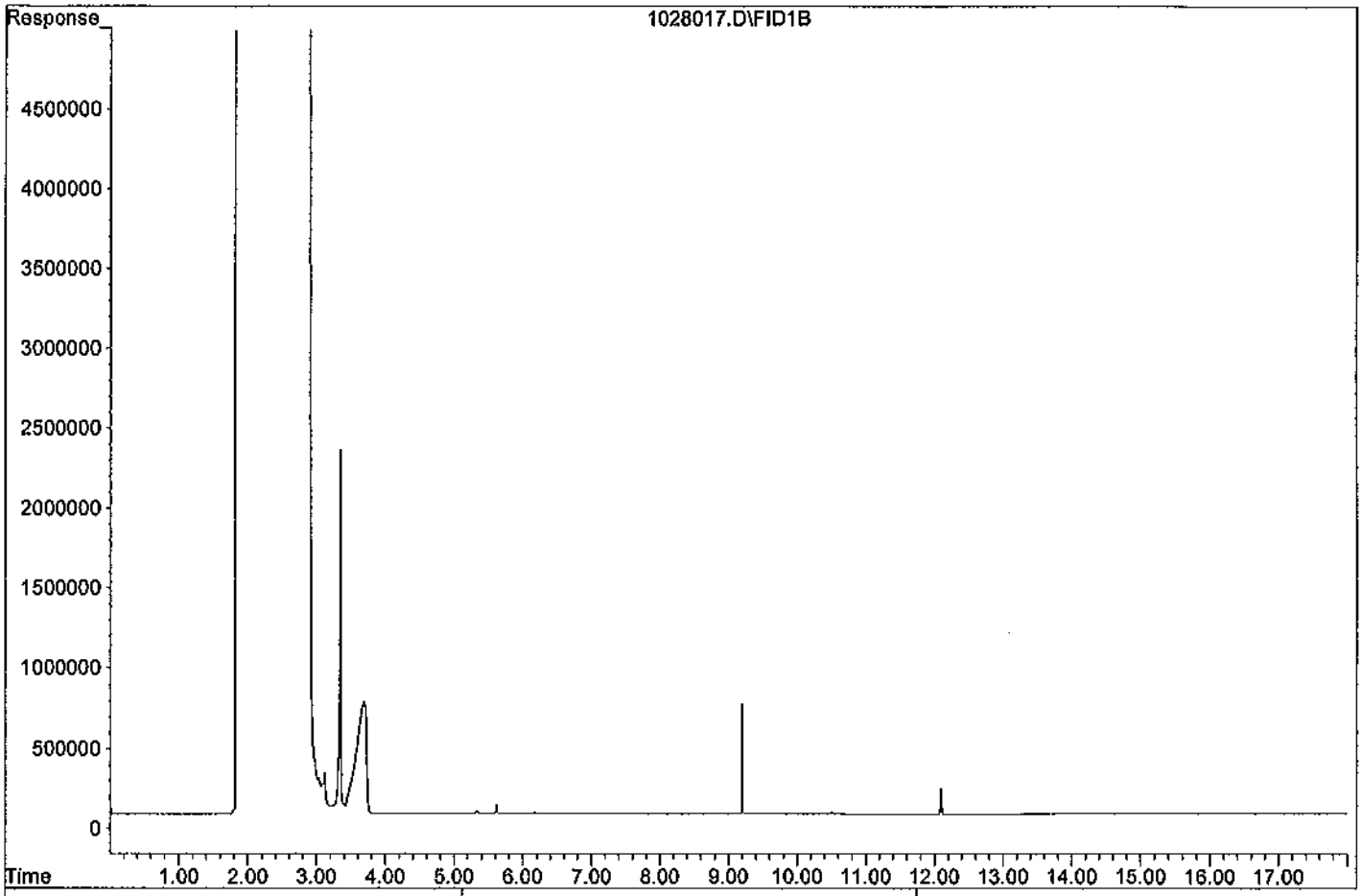
Volume Inj. : 2UL
Signal Phase : DB-5
Signal Info : FID02A

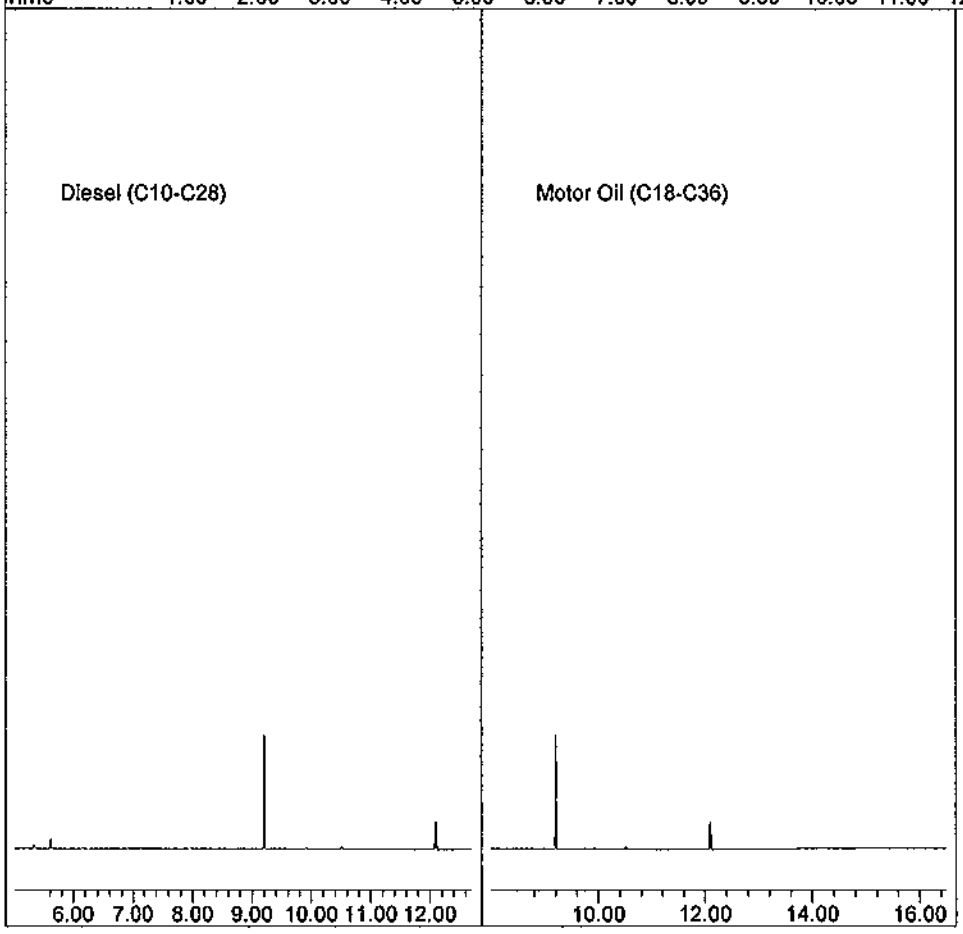
Compound	R.T.	Response	Conc	Units

System Monitoring Compounds

Target Compounds

Data File: G:\APOLLO\DATA\111028\1028017.D
Sample : THC SURR 100/1000





Quantitation Report (Not Reviewed)

Data File : G:\APOLLO\DATA\111028\1028018.D Vial: 18
Acq On : 10-28-11 15:50:20 Operator: LAC
Sample : THC SURR 400/1000 Inst : Apollo
Misc : Mix(C) Multiplr: 1.00
IntFile : events.e
Quant Time: Oct 31 9:01 2011 Quant Results File: TPH1028.RES

Method : G:\APOLLO\DATA\111028\TPH1028.M (Chemstation Integrator)
Title : Diesel
Last Update : Mon Oct 31 10:02:11 2011
Response via : Multiple Level Calibration

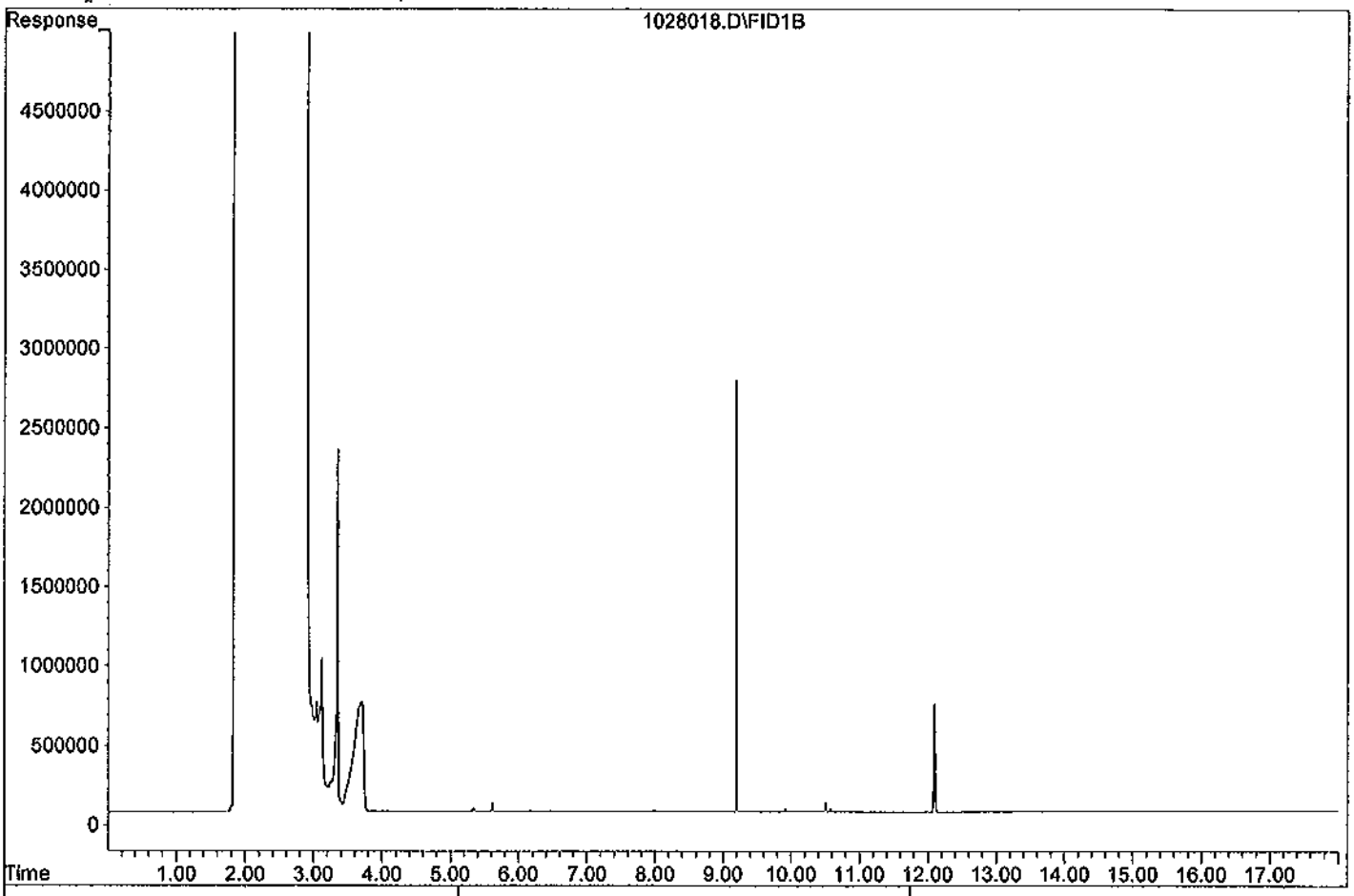
Volume Inj. : 2UL
Signal Phase : DB-5
Signal Info : FID02A

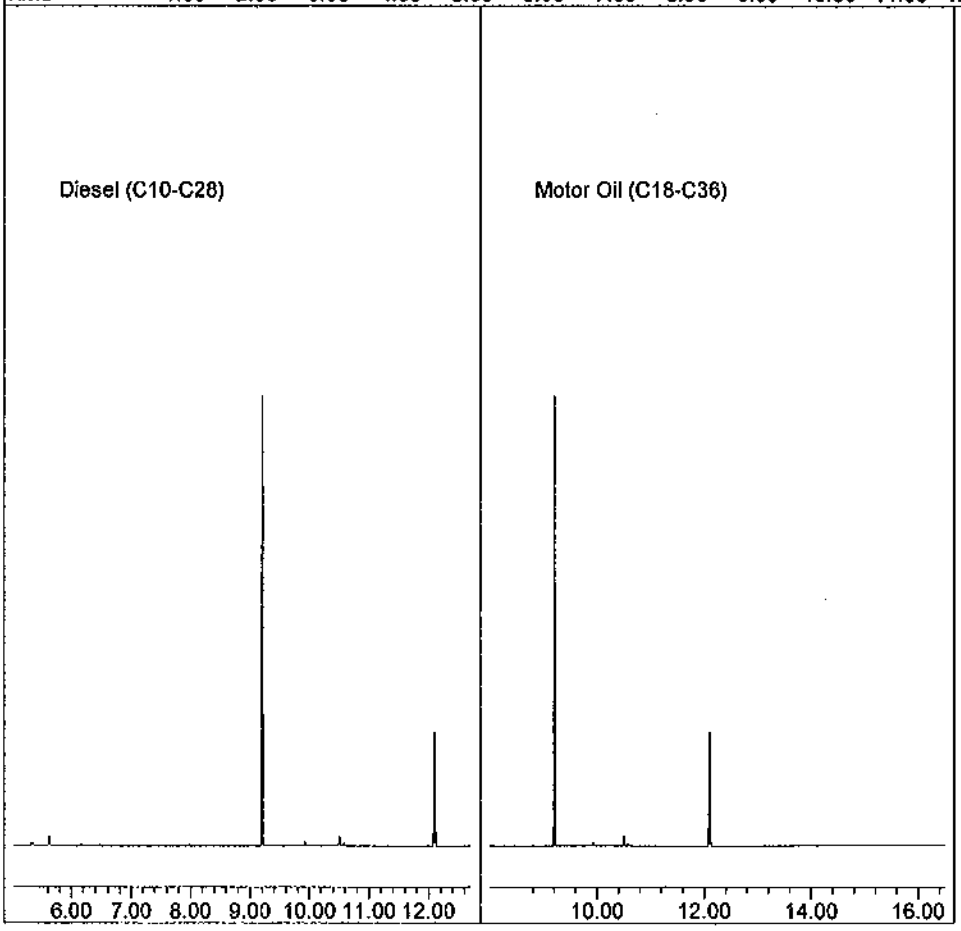
Compound	R.T.	Response	Conc	Units

System Monitoring Compounds

Target Compounds

Data File: G:\APOLLO\DATA\111028\1028018.D
Sample : THC SURR 400/1000





Quantitation Report (Not Reviewed)

Data File : G:\APOLLO\DATA\111028\1028019.D Vial: 19
Acq On : 10-28-11 16:14:52 Operator: LAC
Sample : THC SURR 600/1000 Inst : Apollo
Misc : Mix(C) Multiplr: 1.00
IntFile : events.e
Quant Time: Oct 31 9:01 2011 Quant Results File: TPH1028.RES

Method : G:\APOLLO\DATA\111028\TPH1028.M (Chemstation Integrator)
Title : Diesel
Last Update : Mon Oct 31 10:02:11 2011
Response via : Multiple Level Calibration

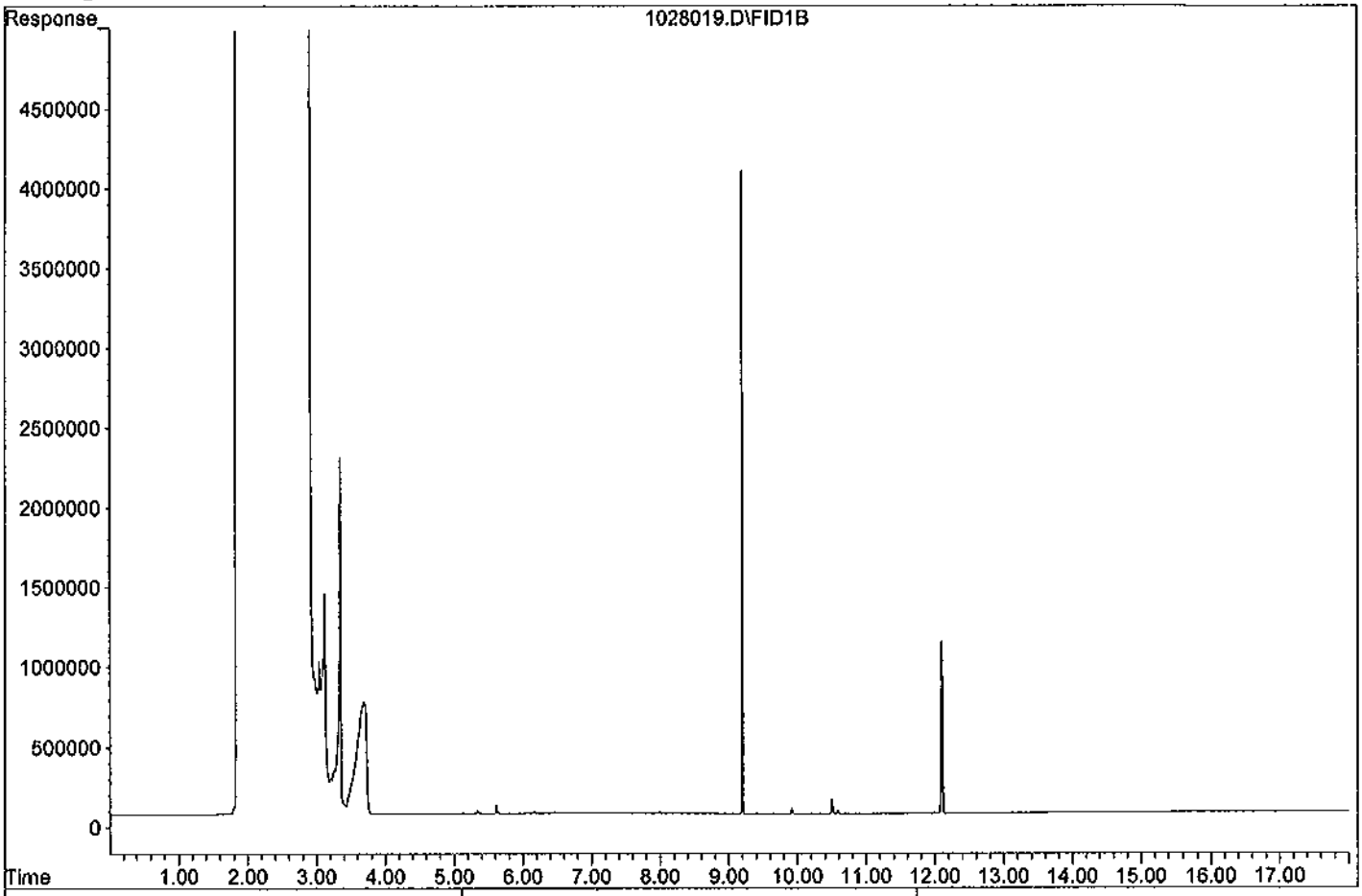
Volume Inj. : 2UL
Signal Phase : DB-5
Signal Info : FID02A

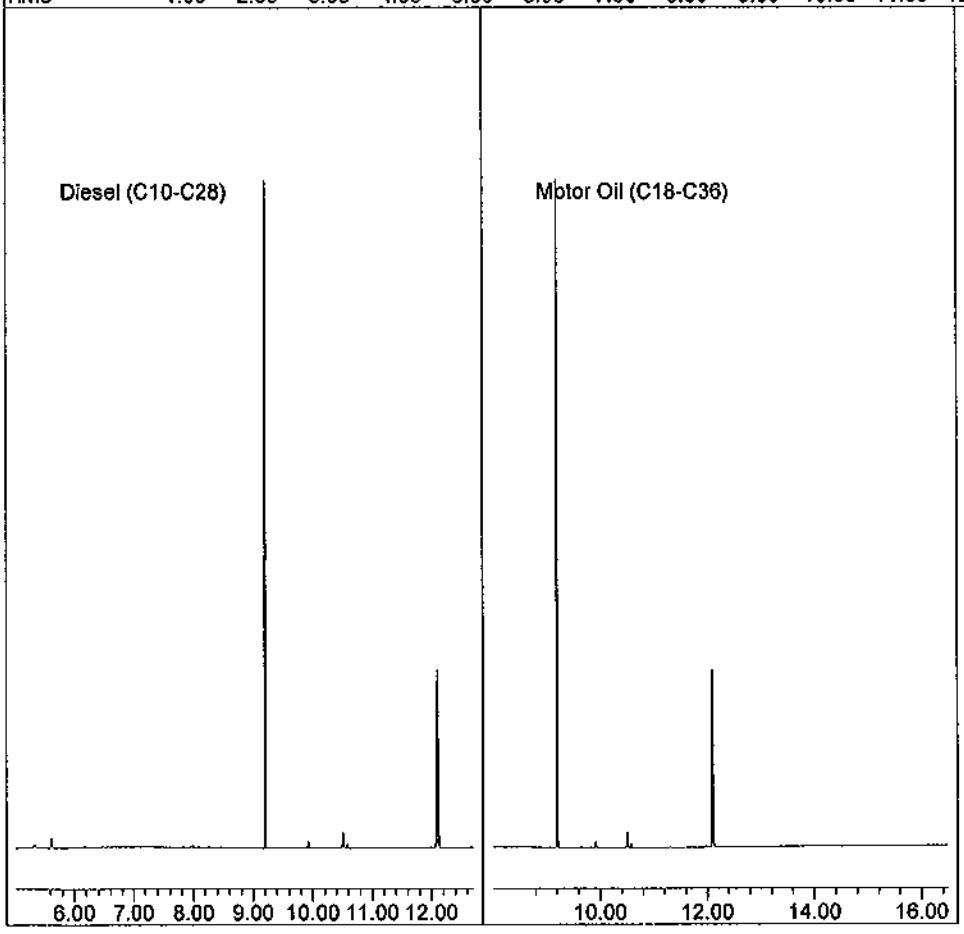
Compound	R.T.	Response	Conc	Units

System Monitoring Compounds

Target Compounds

Data File: G:\APOLLO\DATA\111028\1028019.D
Sample : THC SURR 600/1000





Quantitation Report (Not Reviewed)

Data File : G:\APOLLO\DATA\111028\1028020.D Vial: 20
Acq On : 10-28-11 16:38:57 Operator: LAC
Sample : THC SURR 800/1000 Inst : Apollo
Misc : Mix(C) Multiplr: 1.00
IntFile : events.e
Quant Time: Oct 31 9:01 2011 Quant Results File: TPH1028.RES

Method : G:\APOLLO\DATA\111028\TPH1028.M (Chemstation Integrator)
Title : Diesel
Last Update : Mon Oct 31 10:02:11 2011
Response via : Multiple Level Calibration

Volume Inj. : 2UL
Signal Phase : DB-5
Signal Info : FID02A

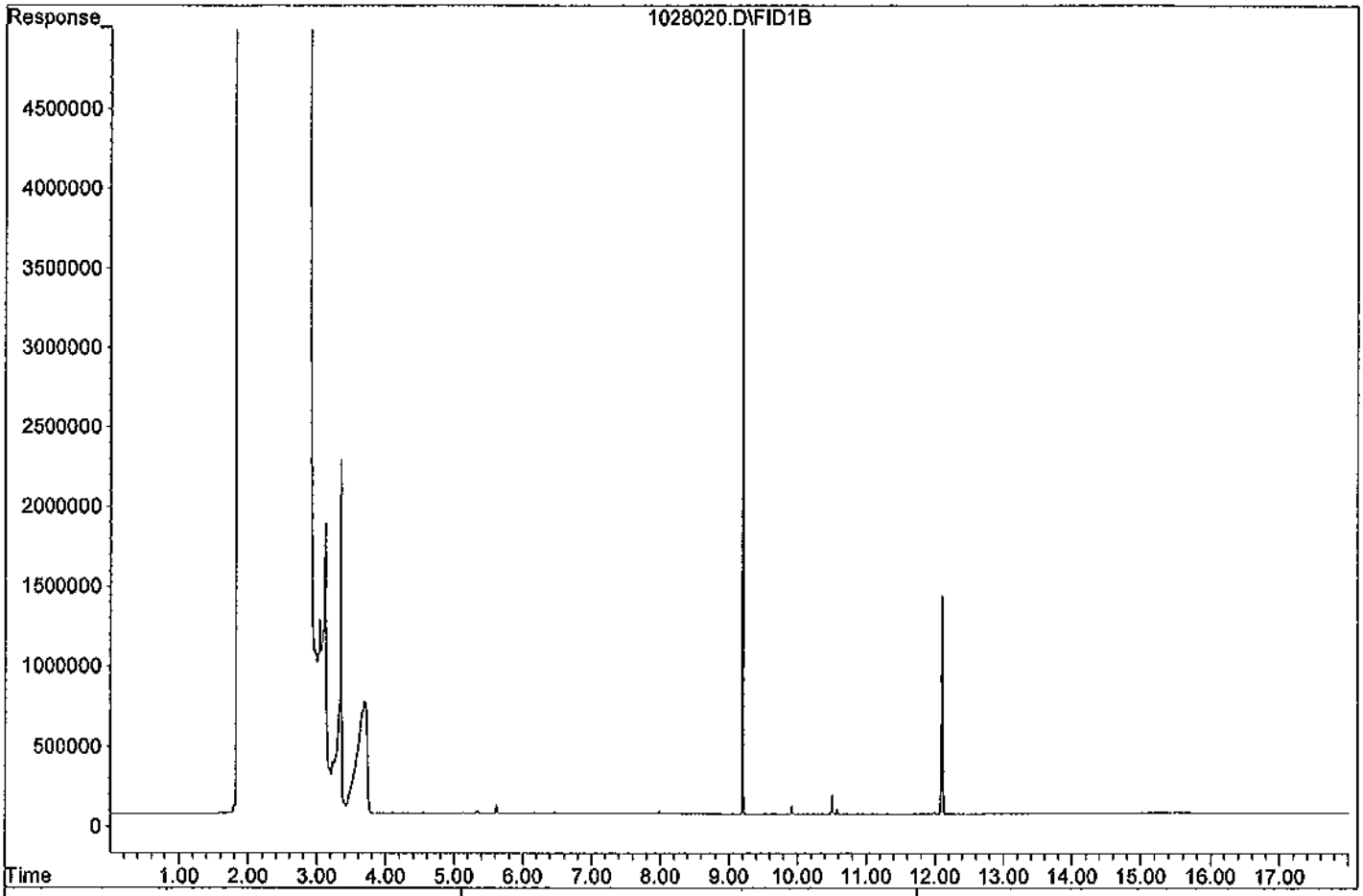
Compound	R.T.	Response	Conc	Units

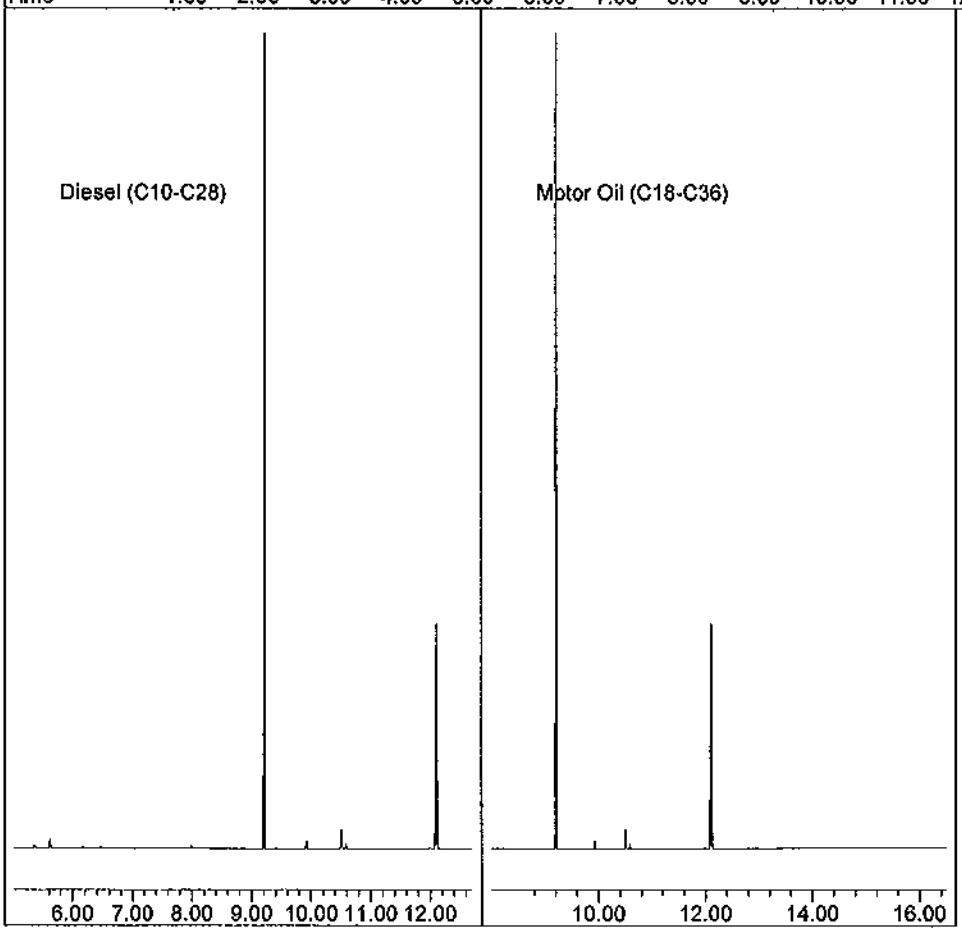
System Monitoring Compounds

Target Compounds

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

Data File: G:\APOLLO\DATA\111028\1028020.D
Sample : THC SURR 800/1000





Quantitation Report (Not Reviewed)

Data File : G:\APOLLO\DATA\111028\1028021.D Vial: 21
Acq On : 10-28-11 17:03:06 Operator: LAC
Sample : THC SURR 1000/1000 Inst : Apollo
Misc : Mix(C) Multiplr: 1.00
IntFile : events.e
Quant Time: Oct 31 9:00 2011 Quant Results File: TPH1028.RES

Method : G:\APOLLO\DATA\111028\TPH1028.M (Chemstation Integrator)
Title : Diesel
Last Update : Mon Oct 31 10:02:11 2011
Response via : Multiple Level Calibration

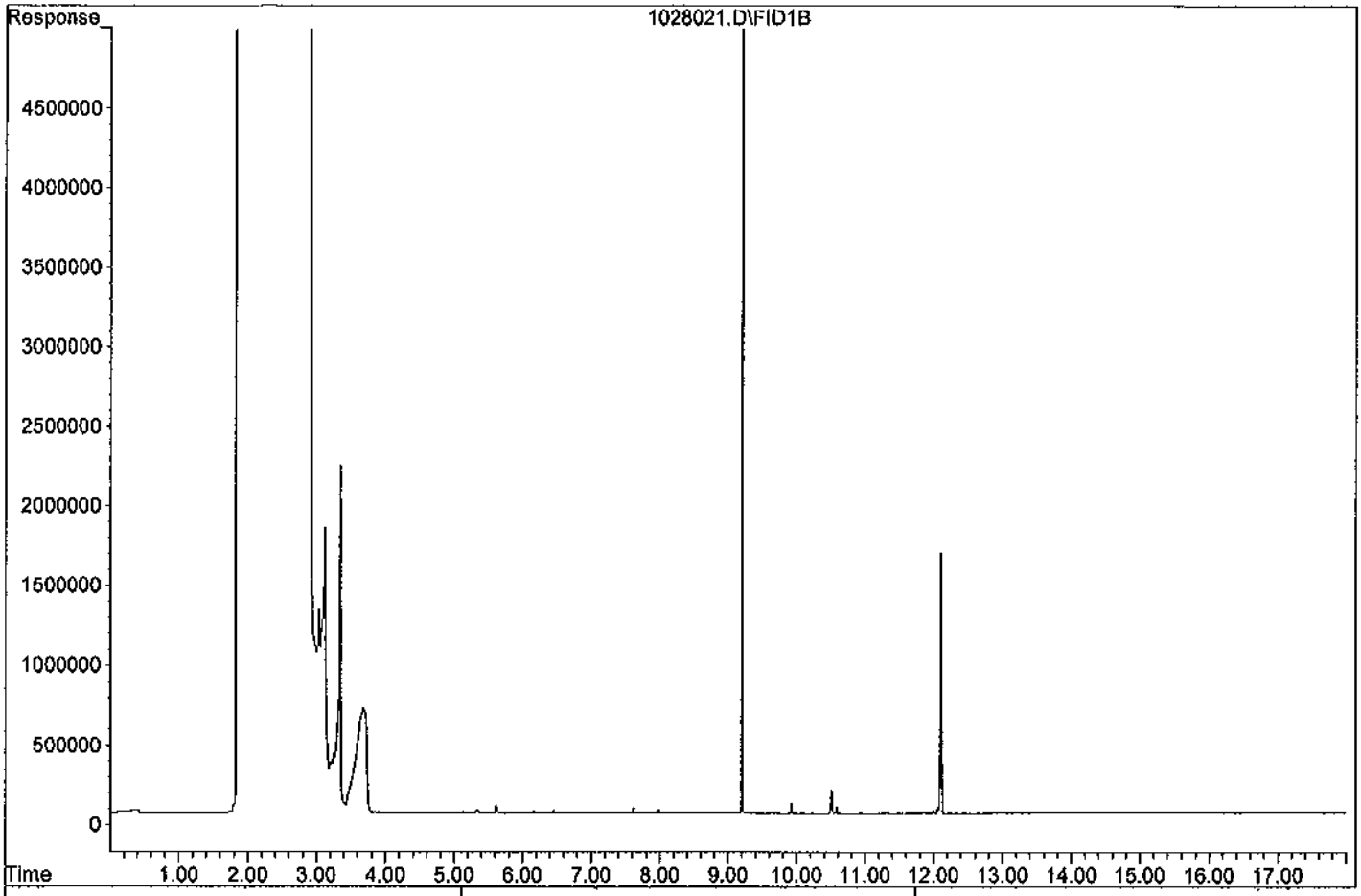
Volume Inj. : 2UL
Signal Phase : DB-5
Signal Info : FID02A

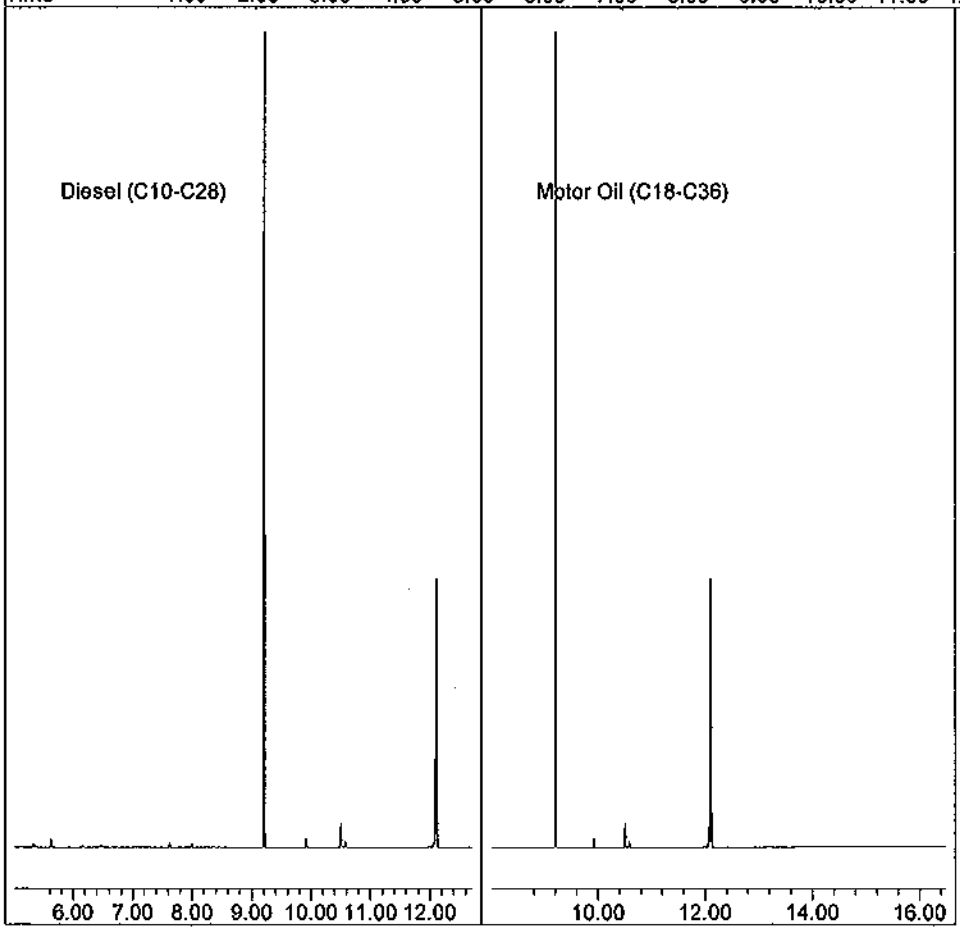
Compound	R.T.	Response	Conc	Units

System Monitoring Compounds

Target Compounds

Data File: G:\APOLLO\DATA\111028\1028021.D
Sample : THC SURR 1000/1000





TPH Extractables
TPH1028

Form 7

Second Source Calibration

Lab Name: APPL, Inc.
Case No:
Matrix:

SDG No: 66102
Date Analyzed: 10/28/11
Instrument: Apollo
Initial Cal. Date: 10/28/11
Data File: 1028015.D

		Compound	MEAN	CCRF	%D		%Drift
1	HATM	Diesel (C10-C28)	420946	437681	4.0	HATM	
2							
3							
4							
5							
6							
7							
8							
9							
10							
11							
12							
13							
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16							
17							
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21							
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27							
28							
29							
30							
31							
32							
33							
34							
35							
36							
37							
38							
39							
40							
		Average			4.0		

Quantitation Report (Not Reviewed)

Data File : G:\APOLLO\DATA\111028\1028015.D Vial: 15
Acq On : 10-28-11 14:37:14 Operator: LAC
Sample : DIESEL 2ND SRC 10/28/11 Inst : Apollo
Misc : Mix(A) Multiplr: 1.00
IntFile : events.e
Quant Time: Oct 28 14:00 2011 Quant Results File: TPH1028.RES

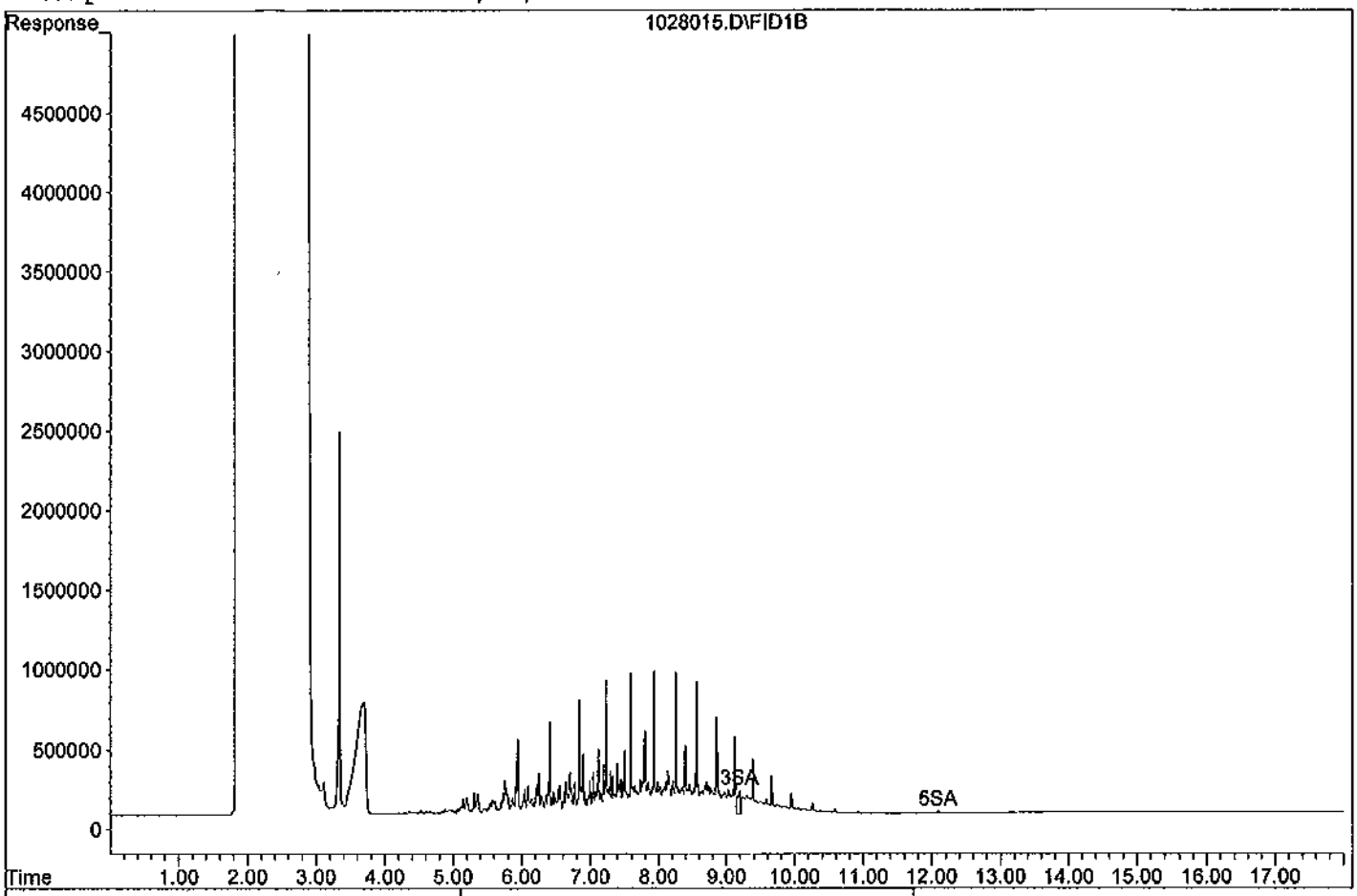
Method : G:\APOLLO\DATA\111028\TPH1028.M (Chemstation Integrator)
Title : Diesel
Last Update : Mon Oct 31 10:02:11 2011
Response via : Multiple Level Calibration

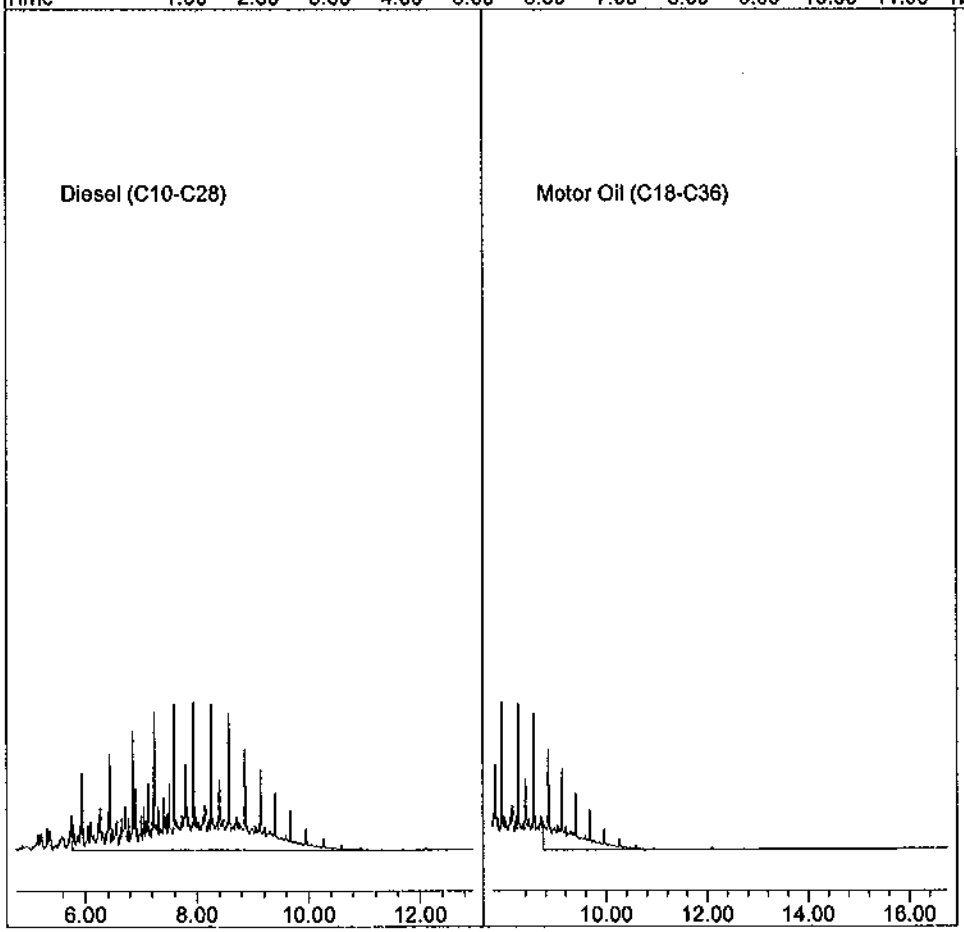
Volume Inj. : 2UL
Signal Phase : DB-5
Signal Info : FID02A

Compound	R.T.	Response	Conc Units
System Monitoring Compounds			
3) SA Not Used(S)	9.20	4372560	4.067 ppb
Surrogate Spike 30.000	Recovery =		13.56%
5) SA Not Used2(S)	12.09	211361	0.427 ppb
Surrogate Spike 30.000	Recovery =		1.42%
Target Compounds			
1) HATM Diesel (C10-C28)	8.86	350144889	415.903 ppb
2) HBTM Motor Oil (C18-C36)	12.25	92370482	254.784 ppb

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

Data File: G:\APOLLO\DATA\111028\1028015.D
Sample : DIESEL 2ND SRC 10/28/11





TPH Extractables
TPH1028

Form 7

Continuing Calibration

Lab Name: APPL, Inc.
Case No:
Matrix:

SDG No: 66102
Date Analyzed: 11/06/11
Instrument: Apollo
Initial Cal. Date: 10/28/11
Data File: 1106003.D

		Compound	MEAN	CCRF	%D		%Drift
1	HATM	Diesel (C10-C28)	420946	359274	15	HATM	
2							
3							
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35							
36							
37							
38							
39							
40							
		Average			15.0		

Quantitation Report (Not Reviewed)

Data File : G:\APOLLO\DATA\111106\1106003.D Vial: 3
Acq On : 11-6-11 16:34:49 Operator: LAC
Sample : DIESEL 400/1000 10/28/11 Inst : Apollo
Misc : Mix(A) Multiplr: 1.00
IntFile : events.e
Quant Time: Nov 7 9:16 2011 Quant Results File: TPH1028.RES

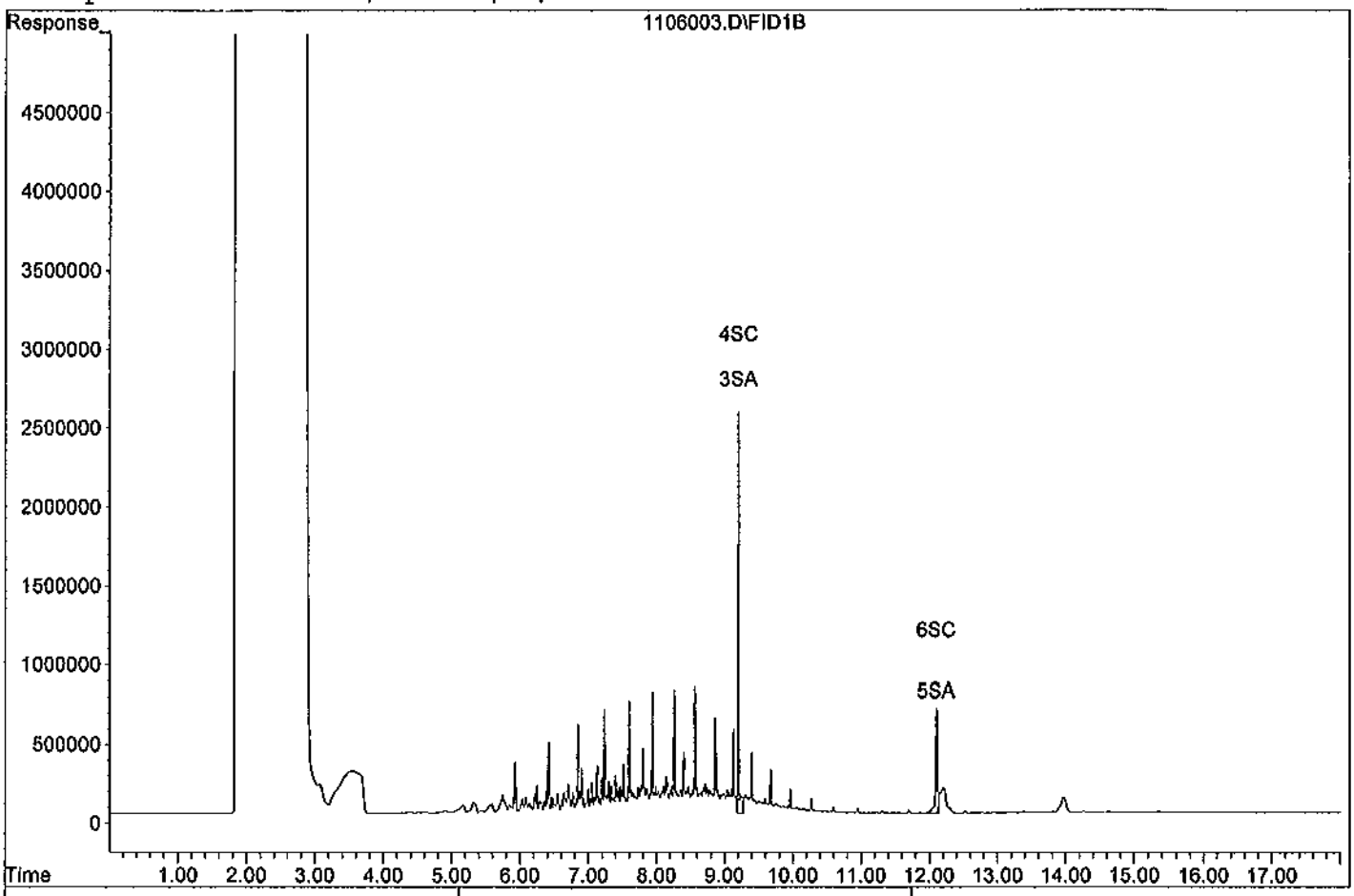
Method : G:\APOLLO\DATA\111028\TPH1028.M (Chemstation Integrator)
Title : Diesel
Last Update : Mon Oct 31 10:02:11 2011
Response via : Multiple Level Calibration

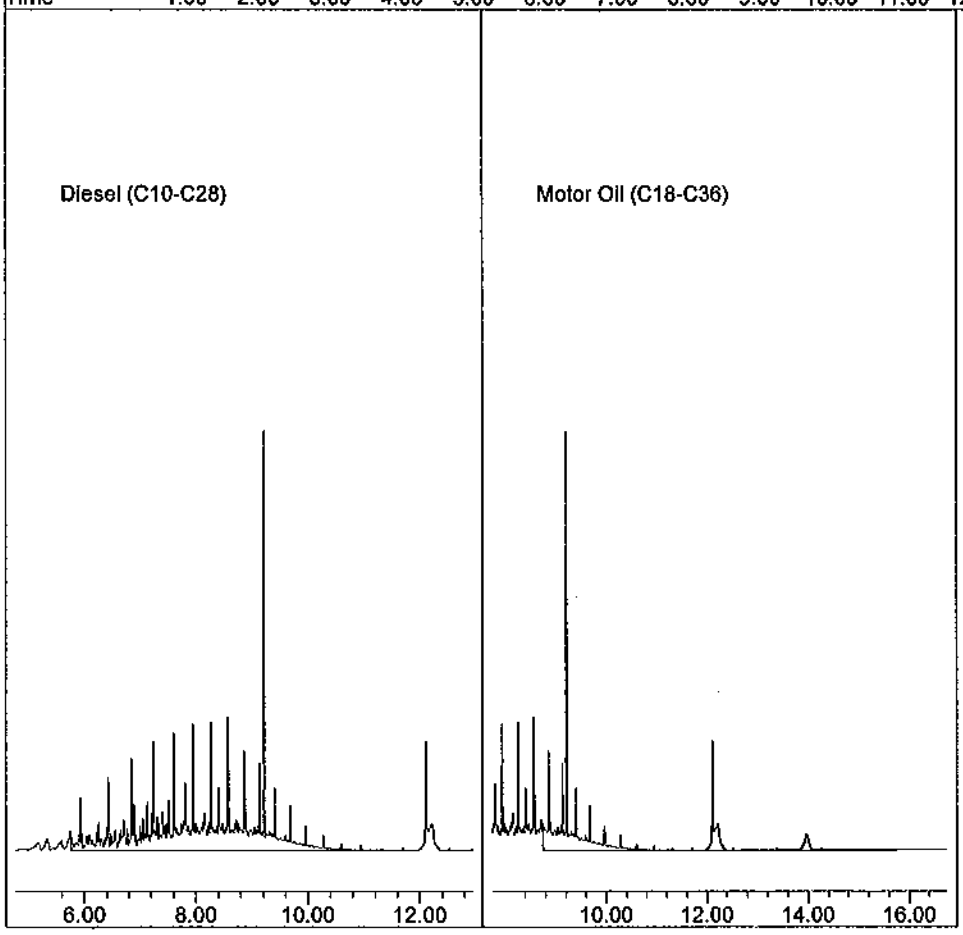
Volume Inj. : 2UL
Signal Phase : DB-5
Signal Info : FID02A

Compound	R.T.	Response	Conc Units
System Monitoring Compounds			
3) SA Not Used(S)	9.21	21346070	19.854 ppb
Surrogate Spike 30.000	Recovery =		66.18%
4) SC Ortho-Terphenyl(S)	9.21	21346070	24.480 ppb
Surrogate Spike 30.000	Recovery =		81.60%
5) SA Not Used2(S)	12.11	11940260	24.100 ppb
Surrogate Spike 30.000	Recovery =		80.33%
6) SC Octacosane(S)	12.11	11940260	25.659 ppb
Surrogate Spike 30.000	Recovery =		85.53%
Target Compounds			
1) HATM Diesel (C10-C28)	8.86	287418867	341.397 ppb
2) HBTM Motor Oil (C18-C36)	12.25	102192727	281.876 ppb

(f)=RT Delta > 1/2 Window 76 (m)=manual int.

Data File: G:\APOLLO\DATA\111106\1106003.D
Sample : DIESEL 400/1000 10/28/11





TPH Extractables
TPH1028

Form 7

Continuing Calibration

Lab Name: APPL, Inc.
Case No:
Matrix:

SDG No: 66102
Date Analyzed: 11/06/11
Instrument: Apollo
Initial Cal. Date: 10/28/11
Data File: 1106017.D

		Compound	MEAN	CCRF	%D		%Drift
1	HATM	Diesel (C10-C28)	420946	369827	12	HATM	
2							
3							
4							
5							
6							
7							
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36							
37							
38							
39							
40							
		Average			12.0		

Quantitation Report (Not Reviewed)

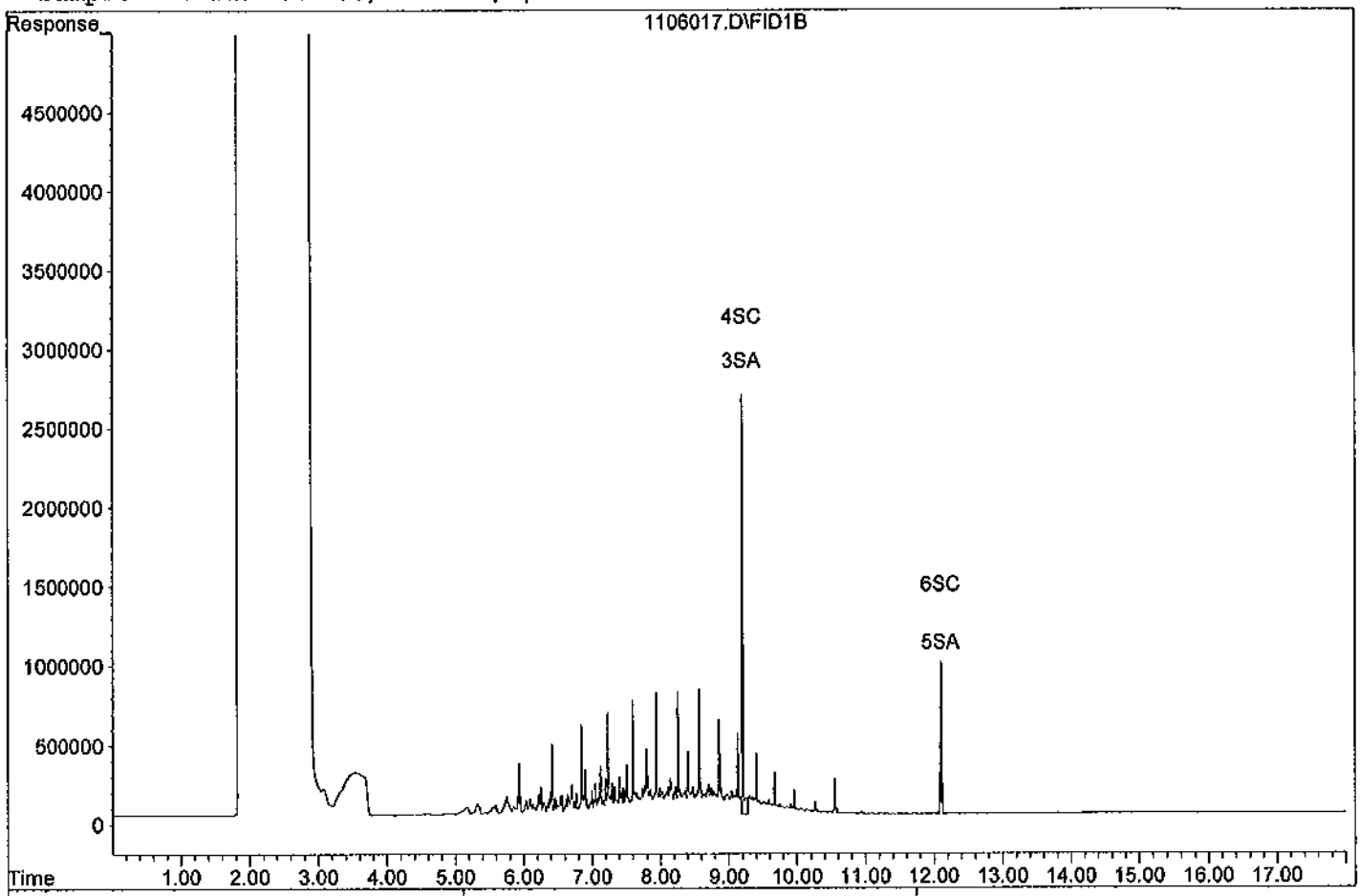
Data File : G:\APOLLO\DATA\111106\1106017.D Vial: 17
Acq On : 11-6-11 22:03:47 Operator: LAC
Sample : DIESEL 400/1000 11/2/11 Inst : Apollo
Misc : Mix(A) Multiplr: 1.00
IntFile : events.e
Quant Time: Nov 7 9:17 2011 Quant Results File: TPH1028.RES

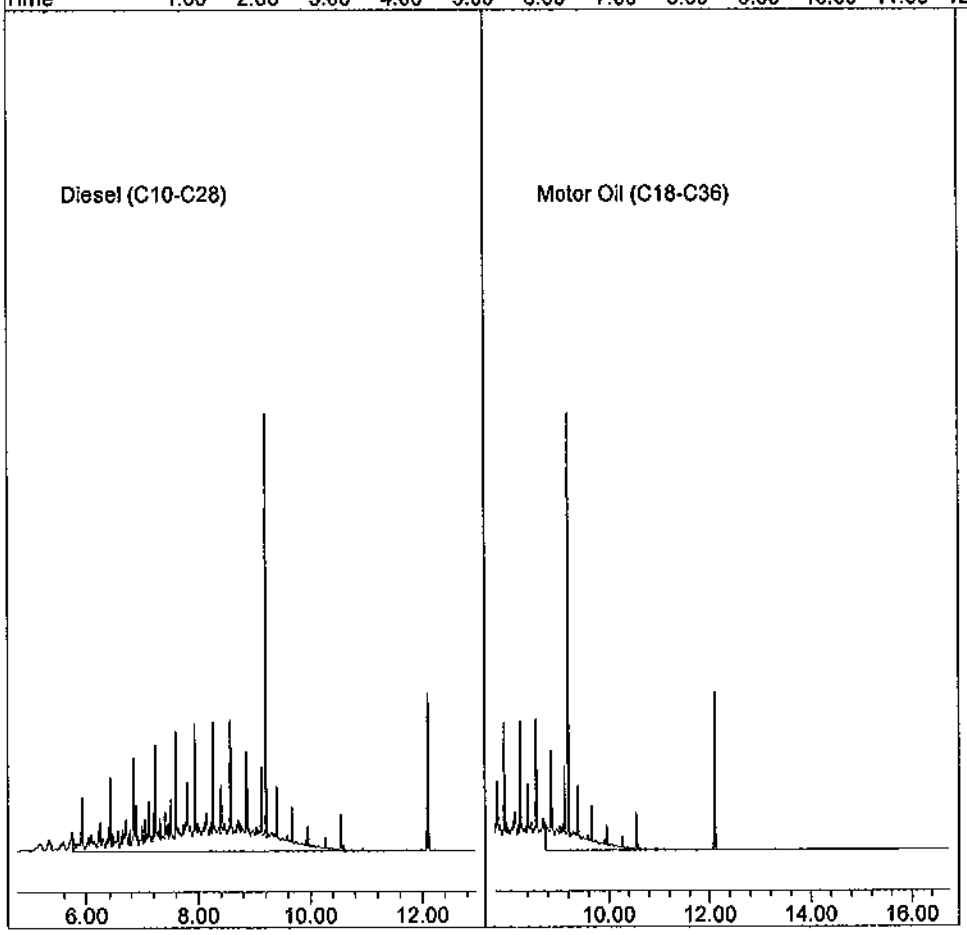
Method : G:\APOLLO\DATA\111028\TPH1028.M (Chemstation Integrator)
Title : Diesel
Last Update : Mon Oct 31 10:02:11 2011
Response via : Multiple Level Calibration

Volume Inj. : 2UL
Signal Phase : DB-5
Signal Info : FID02A

Compound	R.T.	Response	Conc Units
System Monitoring Compounds			
3) SA Not Used(S)	9.21	22135697	20.589 ppb
Surrogate Spike 30.000	Recovery =		68.63%
4) SC Ortho-Terphenyl(S)	9.21	22135697	25.385 ppb
Surrogate Spike 30.000	Recovery =		84.62%
5) SA Not Used2(S)	12.10	13104105	26.450 ppb
Surrogate Spike 30.000	Recovery =		88.17%
6) SC Octacosane(S)	12.10	13104105	28.160 ppb
Surrogate Spike 30.000	Recovery =		93.87%
Target Compounds			
1) HATM Diesel (C10-C28)	8.86	295861444	351.425 ppb
2) HBTM Motor Oil (C18-C36)	12.25	88714324	244.699 ppb

Data File: G:\APOLLO\DATA\111106\1106017.D
Sample : DIESEL 400/1000 11/2/11





TPH Extractables
TPH1028

Form 7

Continuing Calibration

Lab Name: APPL, Inc.
Case No:
Matrix:

SDG No: 66102
Date Analyzed: 11/07/11
Instrument: Apollo
Initial Cal. Date: 10/28/11
Data File: 1106027.D

		Compound	MEAN	CCRF	%D		%Drift
1	HATM	Diesel (C10-C28)	420946	339247	19	HATM	
2							
3							
4							
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		Average			19.0		

Quantitation Report (Not Reviewed)

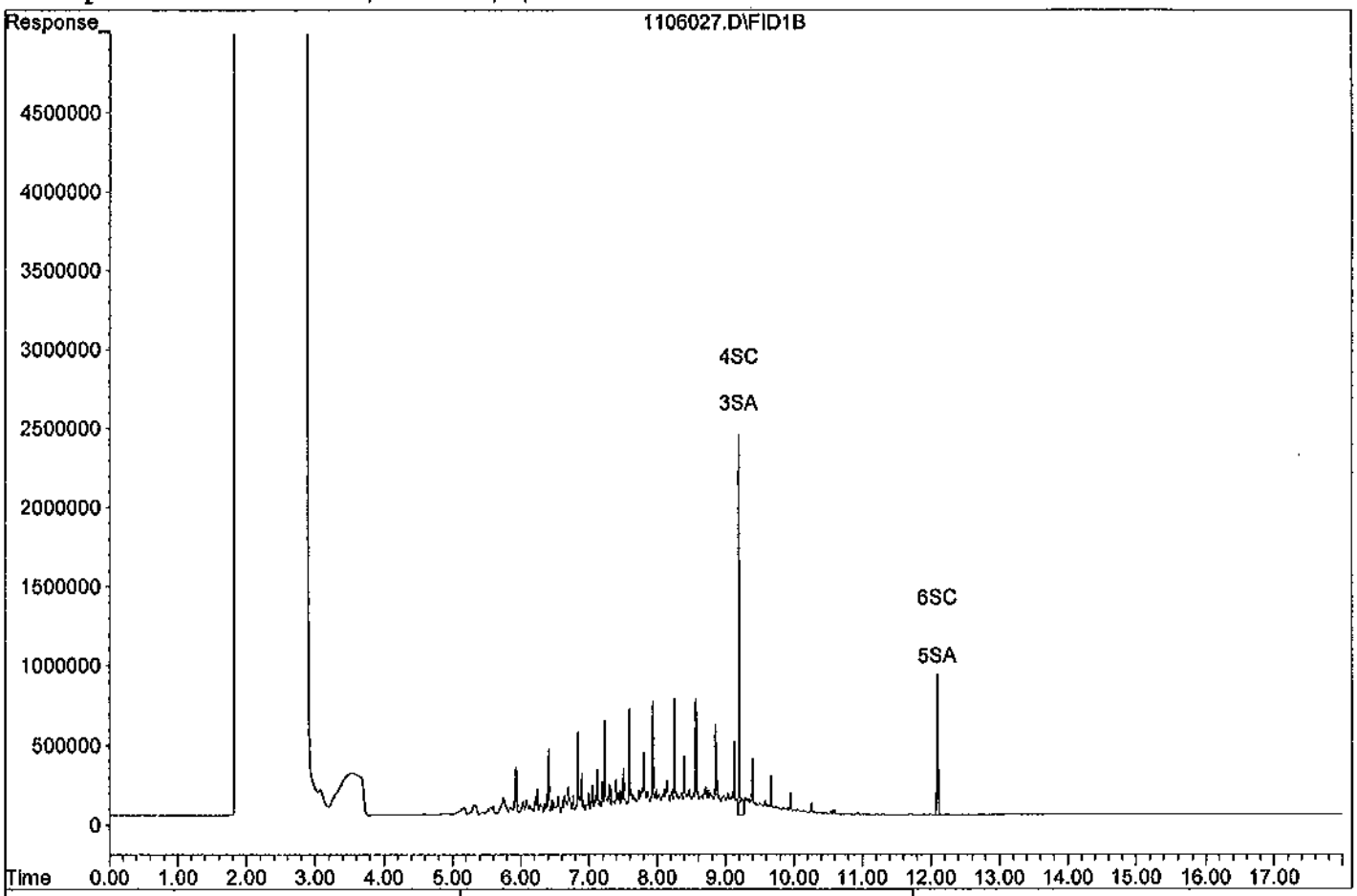
Data File : G:\APOLLO\DATA\111106\1106027.D Vial: 27
Acq On : 11-7-11 1:56:08 Operator: LAC
Sample : DIESEL 400/1000 11/2/11 Inst : Apollo
Misc : Mix(A) Multiplr: 1.00
IntFile : events.e
Quant Time: Nov 7 9:17 2011 Quant Results File: TPH1028.RES

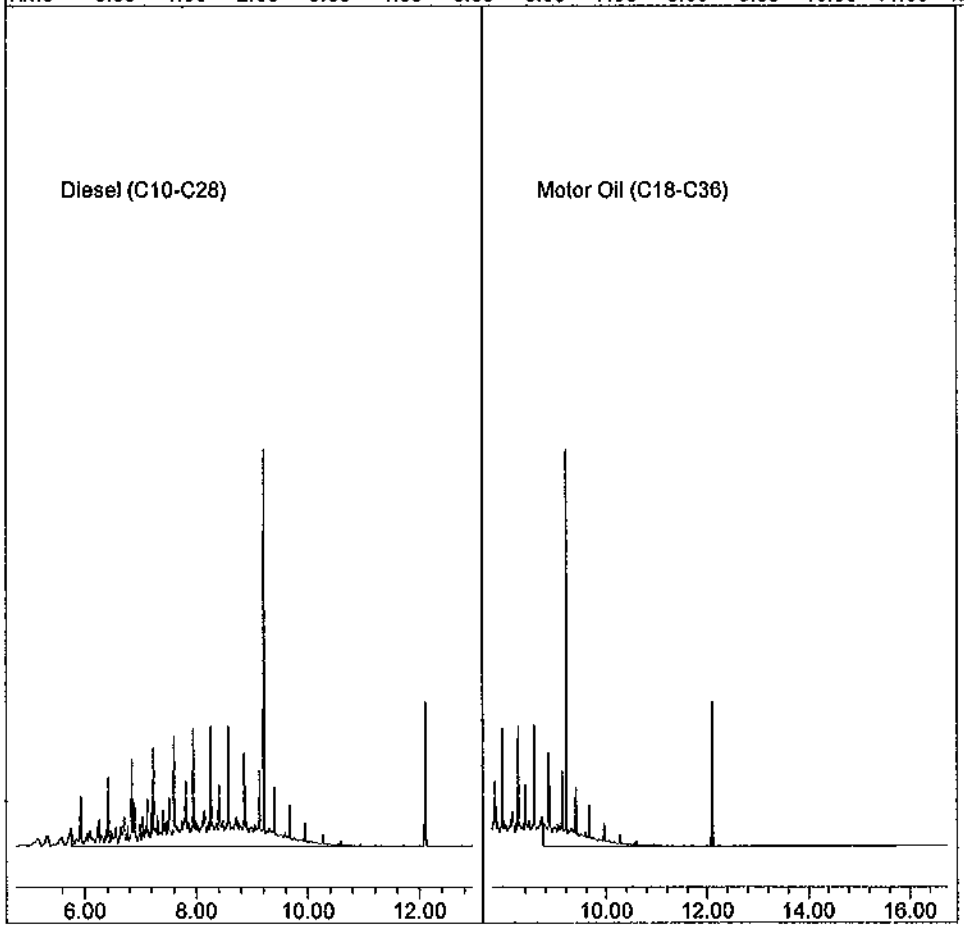
Method : G:\APOLLO\DATA\111028\TPH1028.M (Chemstation Integrator)
Title : Diesel
Last Update : Mon Oct 31 10:02:11 2011
Response via : Multiple Level Calibration

Volume Inj. : 2UL
Signal Phase : DB-5
Signal Info : FID02A

Compound	R.T.	Response	Conc Units
System Monitoring Compounds			
3) SA Not Used(S)	9.21	20565649	19.129 ppb
Surrogate Spike 30.000	Recovery =		63.76%
4) SC Ortho-Terphenyl(S)	9.21	20565649	23.585 ppb
Surrogate Spike 30.000	Recovery =		78.62%
5) SA Not Used2(S)	12.10	11836262	23.890 ppb
Surrogate Spike 30.000	Recovery =		79.63%
6) SC Octacosane(S)	12.10	11836262	25.435 ppb
Surrogate Spike 30.000	Recovery =		84.78%
Target Compounds			
1) HATM Diesel (C10-C28)	8.86	271397500	322.367 ppb
2) HBTM Motor Oil (C18-C36)	12.25	79648905	219.694 ppb

Data File: G:\APOLLO\DATA\111106\1106027.D
Sample : DIESEL 400/1000 11/2/11





TPH Extractables
TPH8S15

Form 6
Initial Calibration

Lab Name: APPL, Inc. SDG No: 66102
Case No: Initial Cal. Date: 11/08/11
Matrix: Instrument: Apollo Initials: LAC

			1115021.D	1115022.D	1115023.D	1115024.D	1115025.D	1115026.D								
		Compound	1	2	3	4	5	6					Avg	%RSD		
1	HATML	Diesel (C10-C28)	613132	243101	243681	243678	244044	245201					305473	49	HATML	1.000
2	HBTM	Motor Oil (C18-C36)	140437	99632	104190	111186	115800	125373					116103	13	HBTM	
3	SA	Not Used(S)	302444	320737	318016	323983	383528	387566					339379	11	SA	
4	SC	Ortho-Terphenyl(S)	356915	320797	300581	304073	324333	307361					319010	6.5	SC	
5	SA	Not Used2(S)		81698	75651	78041	78921	79877					78838	2.8	SA	
6	SC	Octacosane(S)		121445	115156	113245	126484	120297					119325	4.4	SC	
7																
8																
9																
10																
11																
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2.4787102

Quantitation Report (Not Reviewed)

Data File : G:\APOLLO\DATA\111108\1108005.D Vial: 5
Acq On : 11-8-11 15:50:59 Operator: LAC
Sample : DIESEL 100/1000 Inst : Apollo
Misc : Mix(A) Multiplr: 1.00
IntFile : events.e
Quant Time: Nov 16 9:53 2011 Quant Results File: TPH8S15.RES

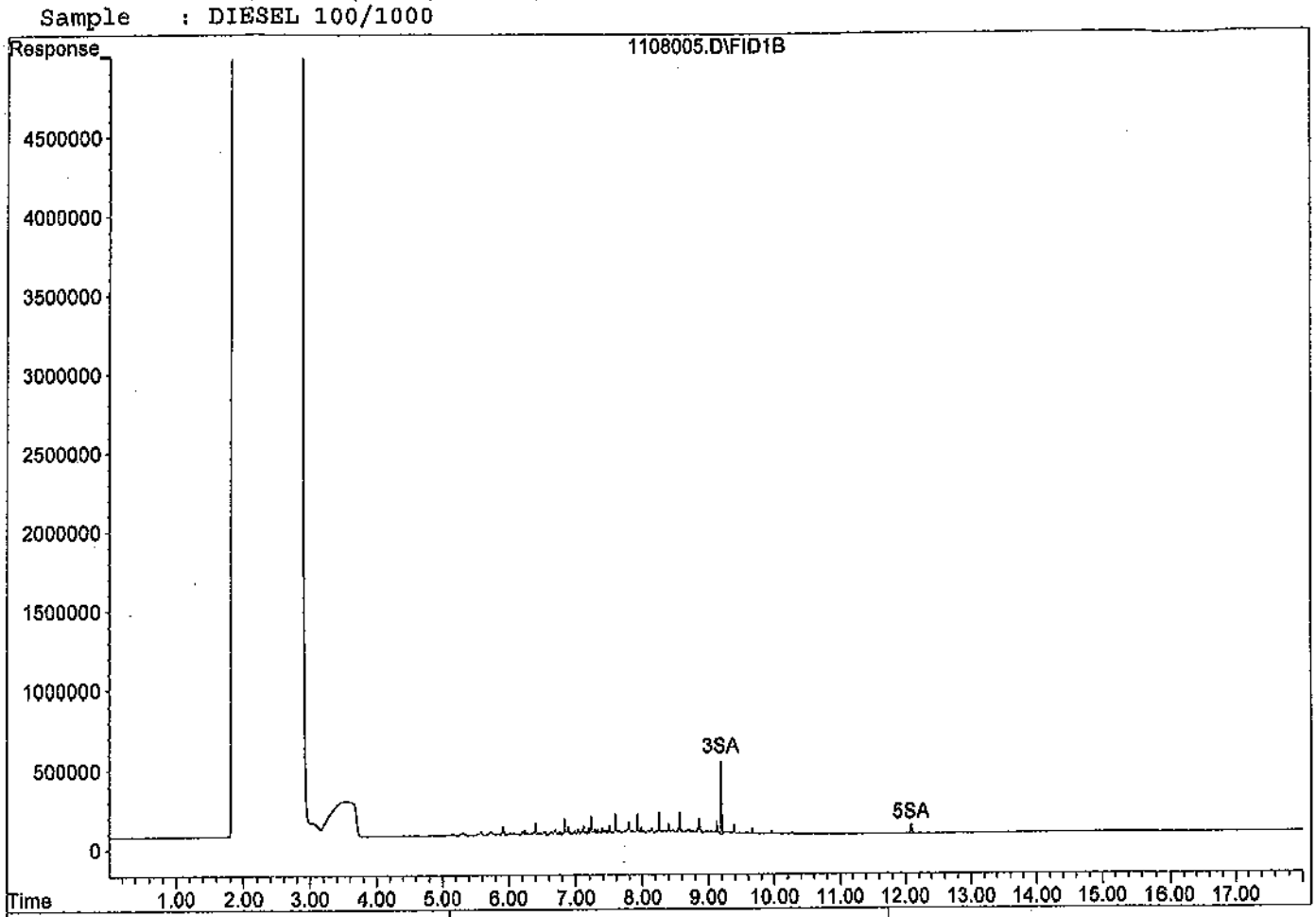
Method : G:\APOLLO\DATA\111115\TPH8S15.M (Chemstation Integrator)
Title : Diesel
Last Update : Thu Nov 17 09:41:49 2011
Response via : Multiple Level Calibration

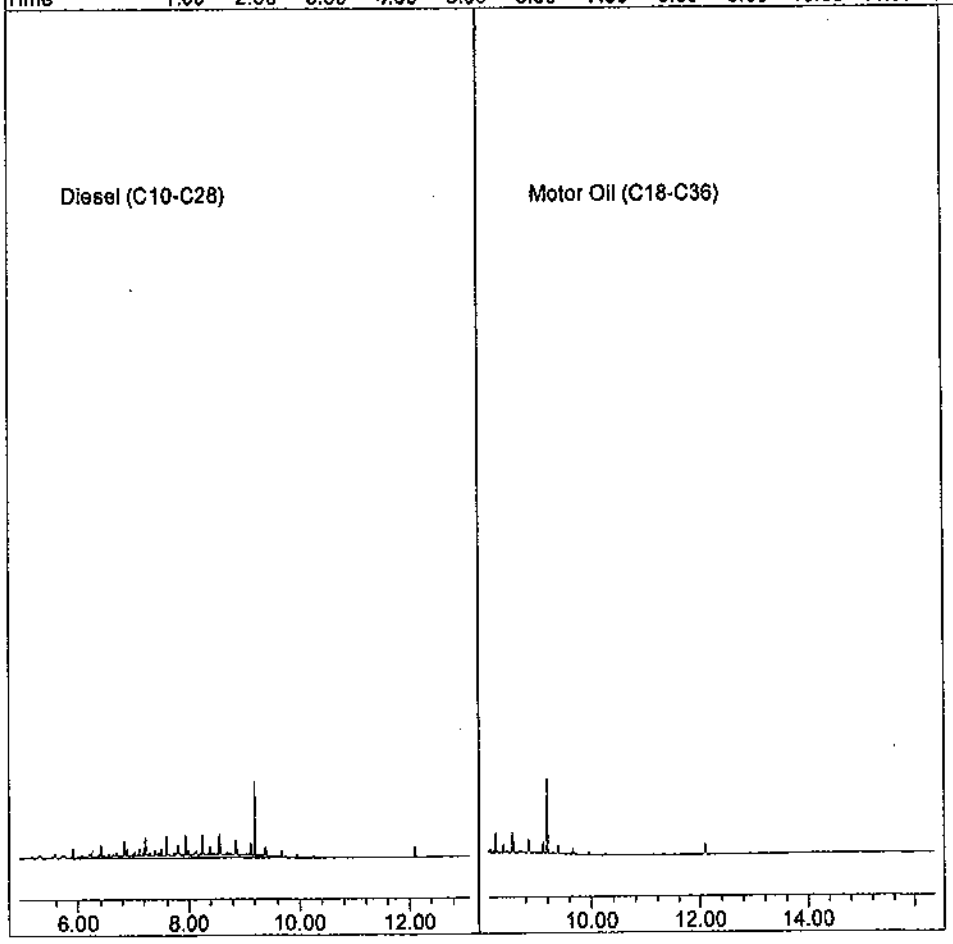
Volume Inj. : 2UL
Signal Phase : DB-5
Signal Info : FID02A

Compound	R.T.	Response	Conc	Units
System Monitoring Compounds				
3) SA Not Used(S)	9.20	3207373	3.154	ppb
Surrogate Spike 30.000	Recovery =		10.51%	
5) SA Not Used2(S)	12.09	816983	2.773	ppb
Surrogate Spike 30.000	Recovery =		9.24%	
Target Compounds				
1) HATM Diesel (C10-C28)	9.01	48620150	767.290	ppb

(f)=RT Delta > 1/2 Window 85 (m)=manual int.

Data File: G:\APOLLO\DATA\111108\1108005.D
Sample : DIESEL 100/1000





Quantitation Report (Not Reviewed)

Data File : G:\APOLLO\DATA\111108\1108006.D Vial: 6
Acq On : 11-8-11 16:14:36 Operator: LAC
Sample : DIESEL 400/1000 Inst : Apollo
Misc : Mix(A) Multiplr: 1.00
IntFile : events.e
Quant Time: Nov 30 11:52 2011 Quant Results File: TPH8S15.RES

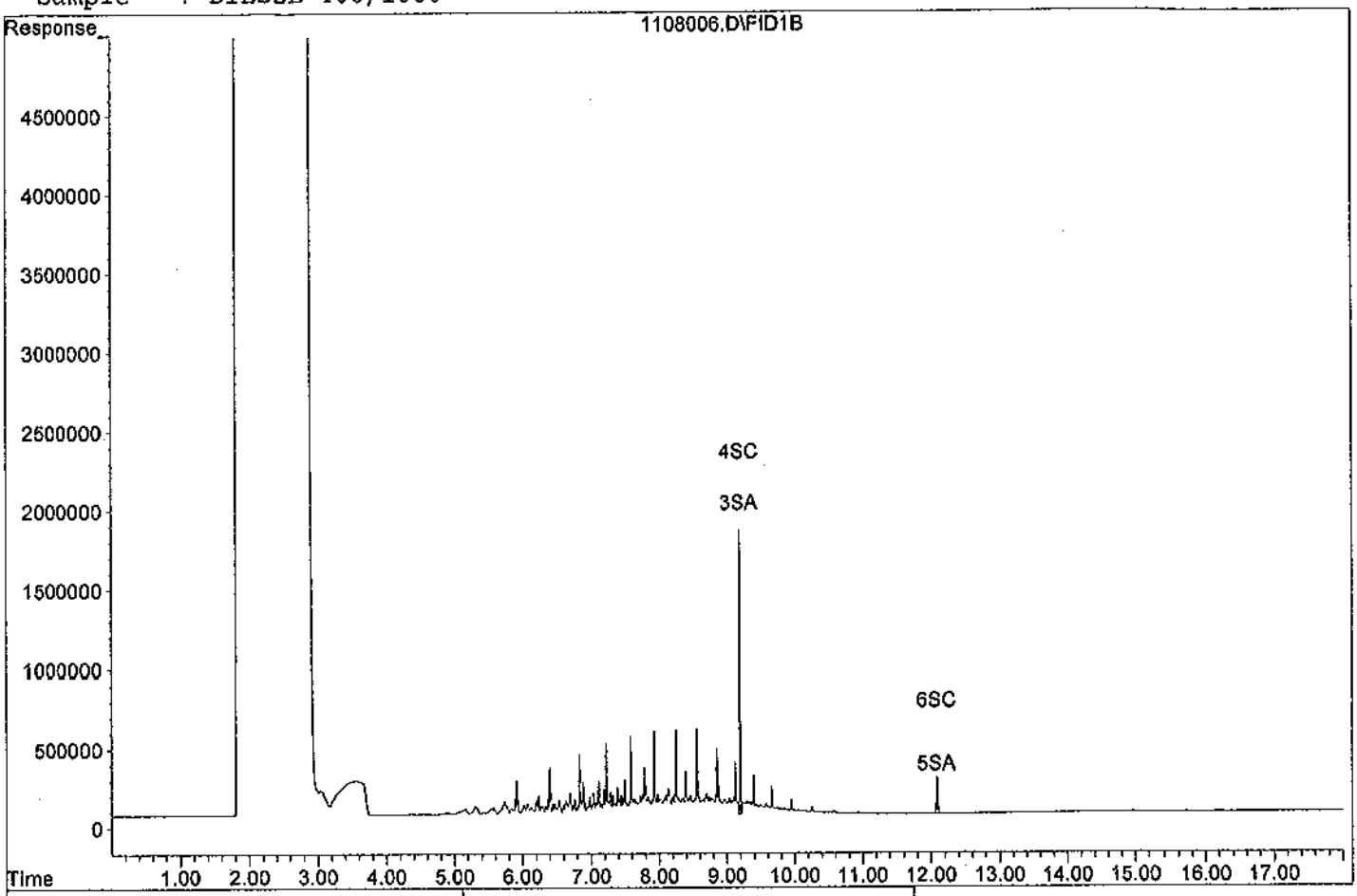
Method : G:\APOLLO\DATA\111108\TPH8S15.M (Chemstation Integrator)
Title : Diesel
Last Update : Wed Nov 30 11:52:46 2011
Response via : Multiple Level Calibration

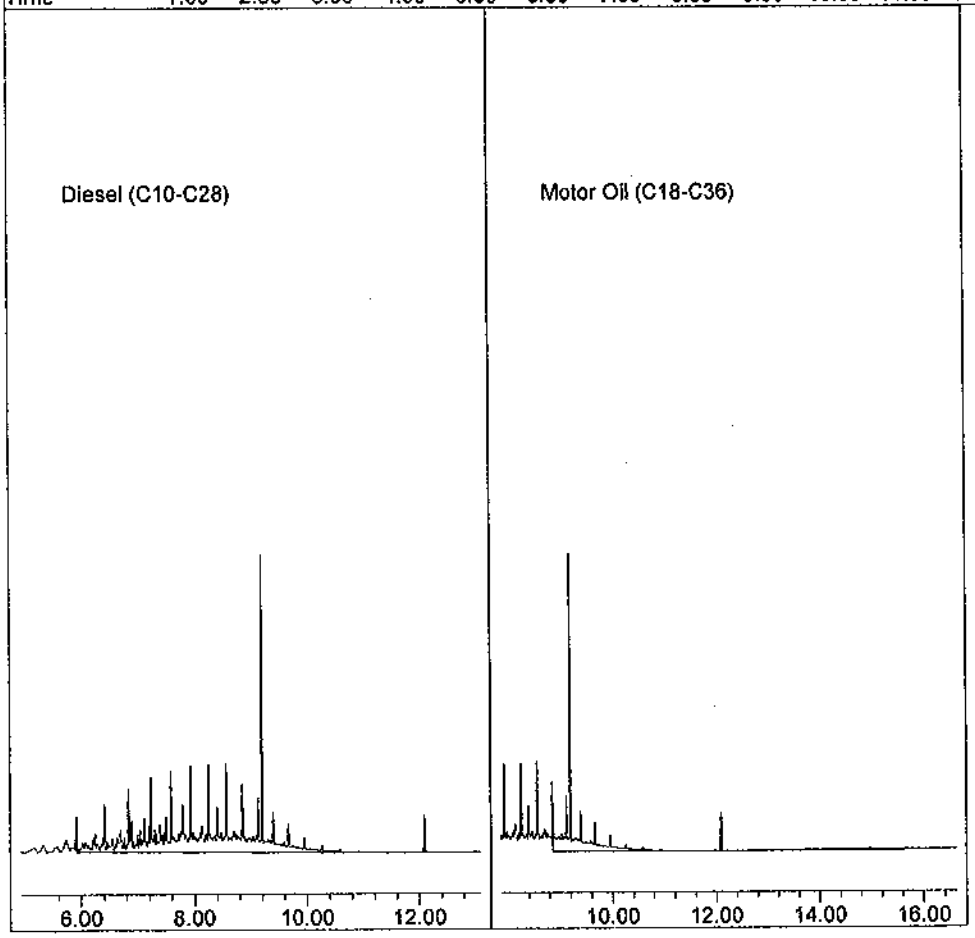
Volume Inj. : 2UL
Signal Phase : DB-5
Signal Info : FID02A

Compound	R.T.	Response	Conc Units
System Monitoring Compounds			
3) SA Not Used(S)	9.20	12720627	18.741 ppb
Surrogate Spike 30.000	Recovery =		62.47%
4) SC Ortho-Terphenyl(S)	9.20	12720627	19.938 ppb
Surrogate Spike 30.000	Recovery =		66.46%
5) SA Not Used2(S)	12.09	3026041	19.192 ppb
Surrogate Spike 30.000	Recovery =		63.97%
6) SC Octacosane(S)	12.09	3026041	12.960 ppb
Surrogate Spike 30.000	Recovery =		43.20%
Target Compounds			
1) HATM Diesel (C10-C28)	9.01	194945056	395.862 ppb
2) HBTM Motor Oil (C18-C36)	12.24	62241854	268.046 ppb

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

Data File: G:\APOLLO\DATA\111108\1108006.D
Sample : DIESEL 400/1000





Quantitation Report (Not Reviewed)

Data File : G:\APOLLO\DATA\111108\1108007.D Vial: 7
Acq On : 11-8-11 16:38:14 Operator: LAC
Sample : DIESEL 600/1000 Inst : Apollo
Misc : Mix(A) Multiplr: 1.00
IntFile : events.e
Quant Time: Nov 30 11:53 2011 Quant Results File: TPH8S15.RES

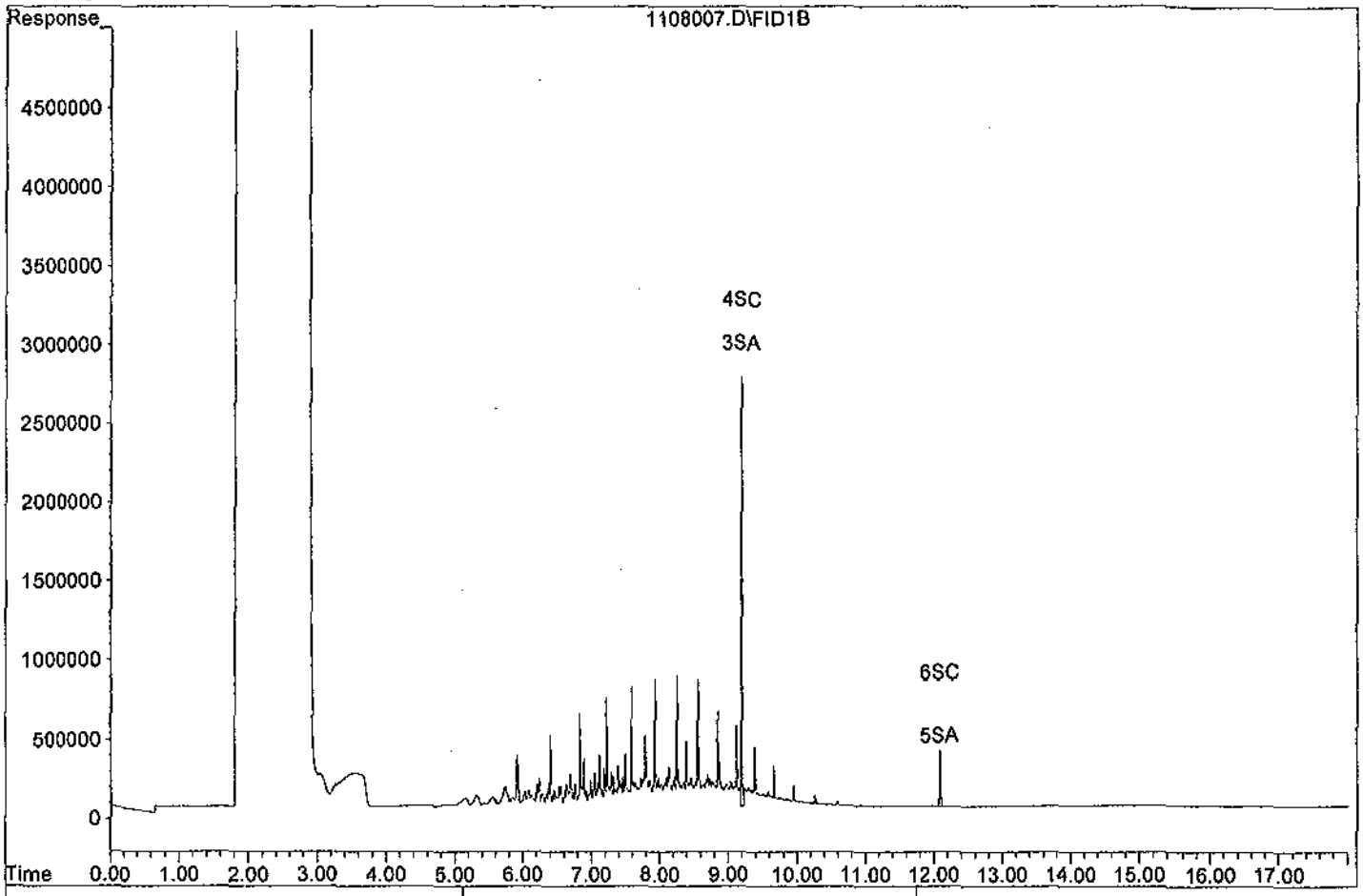
Method : G:\APOLLO\DATA\111108\TPH8S15.M (Chemstation Integrator)
Title : Diesel
Last Update : Wed Nov 30 11:52:46 2011
Response via : Multiple Level Calibration

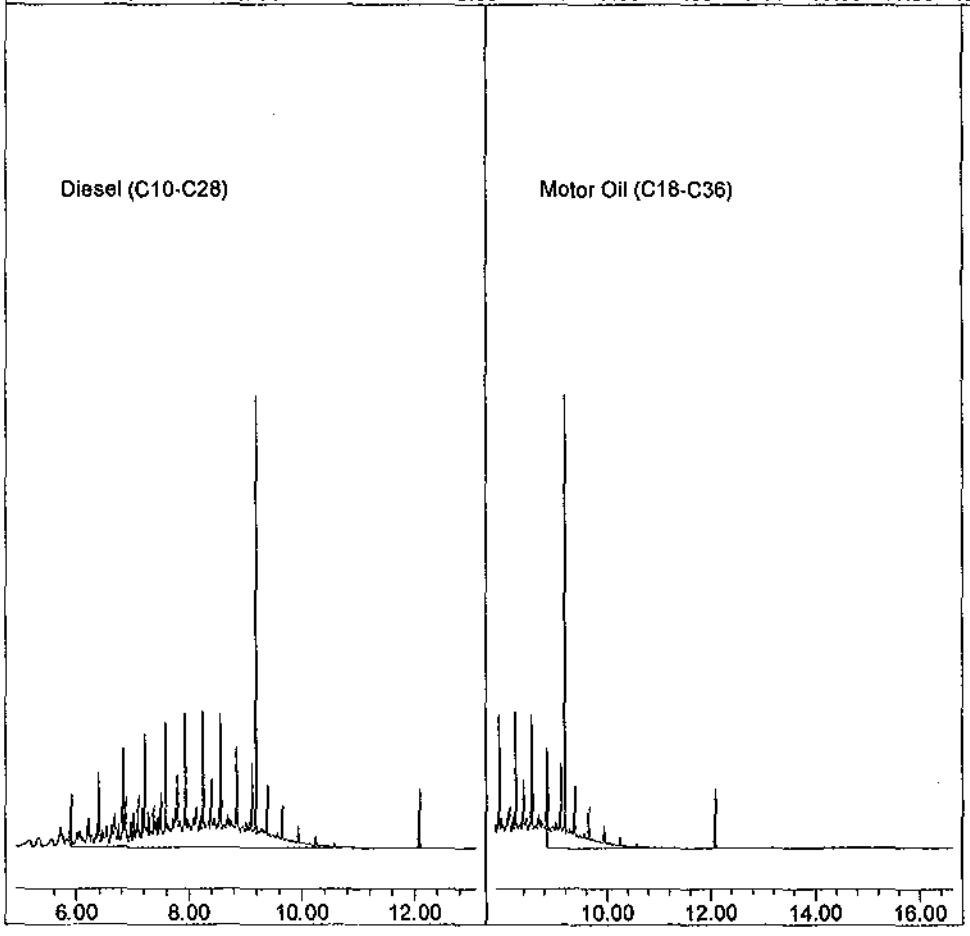
Volume Inj. : 2UL
Signal Phase : DB-5
Signal Info : FID02A

Compound	R.T.	Response	Conc	Units
System Monitoring Compounds				
3) SA Not Used(S)	9.20	19438997	28.639	ppb
Surrogate Spike 30.000	Recovery =		95.46%	
4) SC Ortho-Terphenyl(S)	9.20	19438997	30.468	ppb
Surrogate Spike 30.000	Recovery =		101.56%	
5) SA Not Used2(S)	12.09	4682445	29.697	ppb
Surrogate Spike 30.000	Recovery =		98.99%	
6) SC Octacosane(S)	12.09	4682445	20.054	ppb
Surrogate Spike 30.000	Recovery =		66.85%	
Target Compounds				
1) HATM Diesel (C10-C28)	9.01	292413883	596.788	ppb
2) HBTM Motor Oil (C18-C36)	12.24	77206924	332.493	ppb

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

Data File: G:\APOLLO\DATA\111108\1108007.D
Sample : DIESEL 600/1000





Quantitation Report (Not Reviewed)

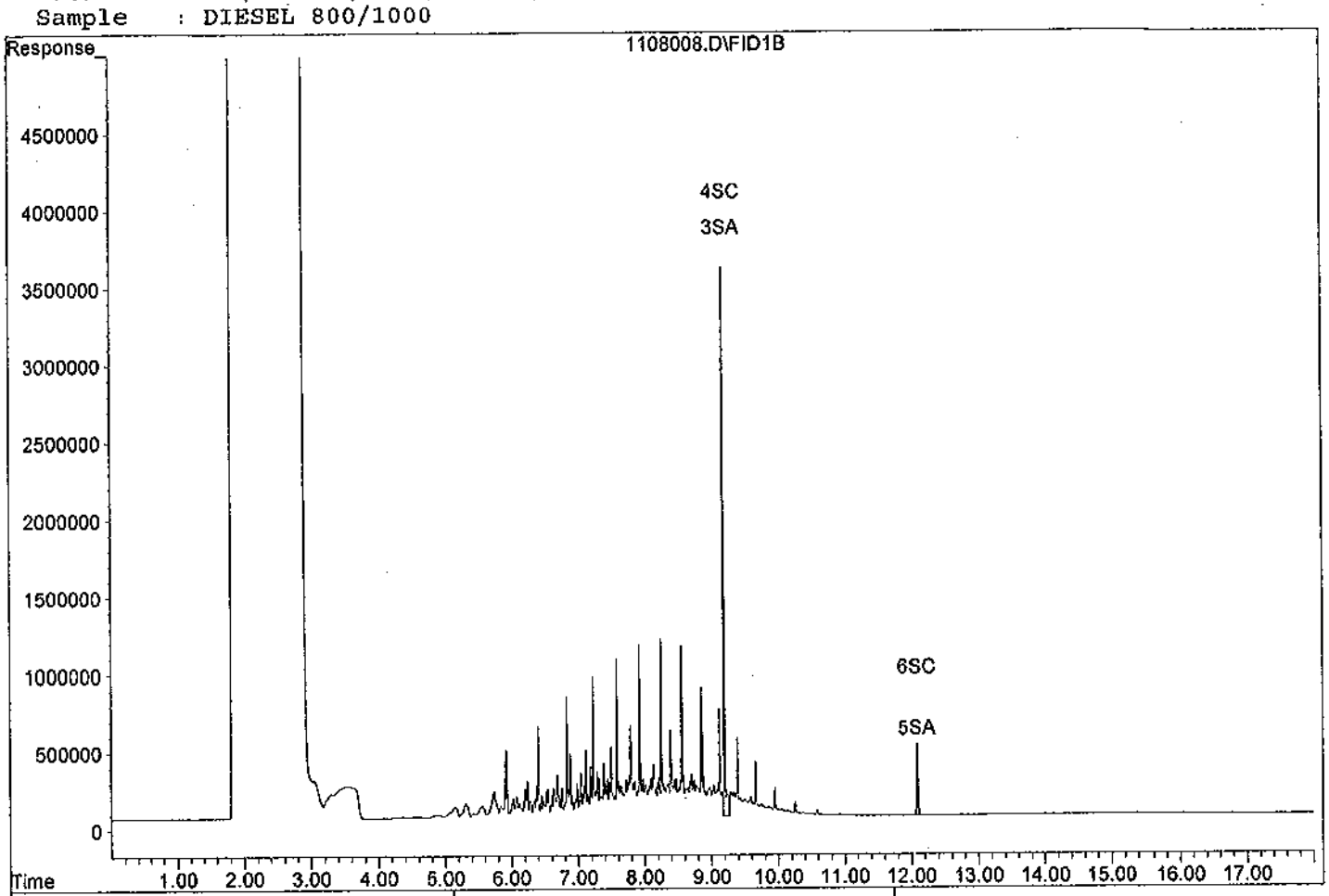
Data File : G:\APOLLO\DATA\111108\1108008.D Vial: 8
Acq On : 11-8-11 17:01:53 Operator: LAC
Sample : DIESEL 800/1000 Inst : Apollo
Misc : Mix(A) Multiplr: 1.00
IntFile : events.e
Quant Time: Nov 30 11:53 2011 Quant Results File: TPH8S15.RES

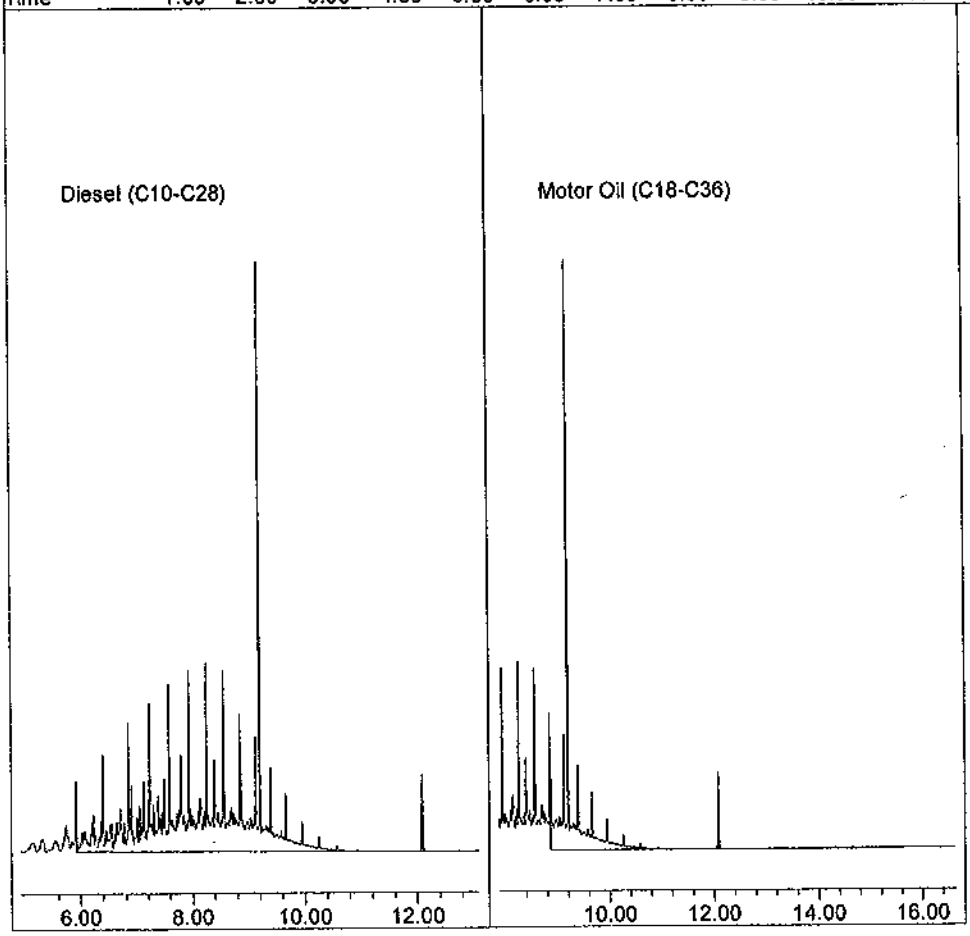
Method : G:\APOLLO\DATA\111108\TPH8S15.M (Chemstation Integrator)
Title : Diesel
Last Update : Wed Nov 30 11:52:46 2011
Response via : Multiple Level Calibration

Volume Inj. : 2UL
Signal Phase : DB-5
Signal Info : FID02A

Compound	R.T.	Response	Conc	Units
System Monitoring Compounds				
3) SA Not Used(S)	9.20	30682231	45.203	ppb
Surrogate Spike 30.000	Recovery =		150.68%	
4) SC Ortho-Terphenyl(S)	9.20	30682231	48.090	ppb
Surrogate Spike 30.000	Recovery =		160.30%	
5) SA Not Used2(S)	12.09	6313667	40.042	ppb
Surrogate Spike 30.000	Recovery =		133.47%	
6) SC Octacosane(S)	12.09	6313667	27.041	ppb
Surrogate Spike 30.000	Recovery =		90.14%	
Target Compounds				
1) HATM Diesel (C10-C28)	9.01	390470225	798.924	ppb
2) HBTM Motor Oil (C18-C36)	12.24	99027136	426.462	ppb

Data File: G:\APOLLO\DATA\111108\1108008.D
Sample : DIESEL 800/1000





Quantitation Report (Not Reviewed)

Data File : G:\APOLLO\DATA\111108\1108009.D Vial: 9
Acq On : 11-8-11 17:25:32 Operator: LAC
Sample : DIESEL 1000/1000 Inst : Apollo
Misc : Mix(A) Multiplr: 1.00
IntFile : events.e
Quant Time: Nov 30 11:53 2011 Quant Results File: TPH8S15.RES

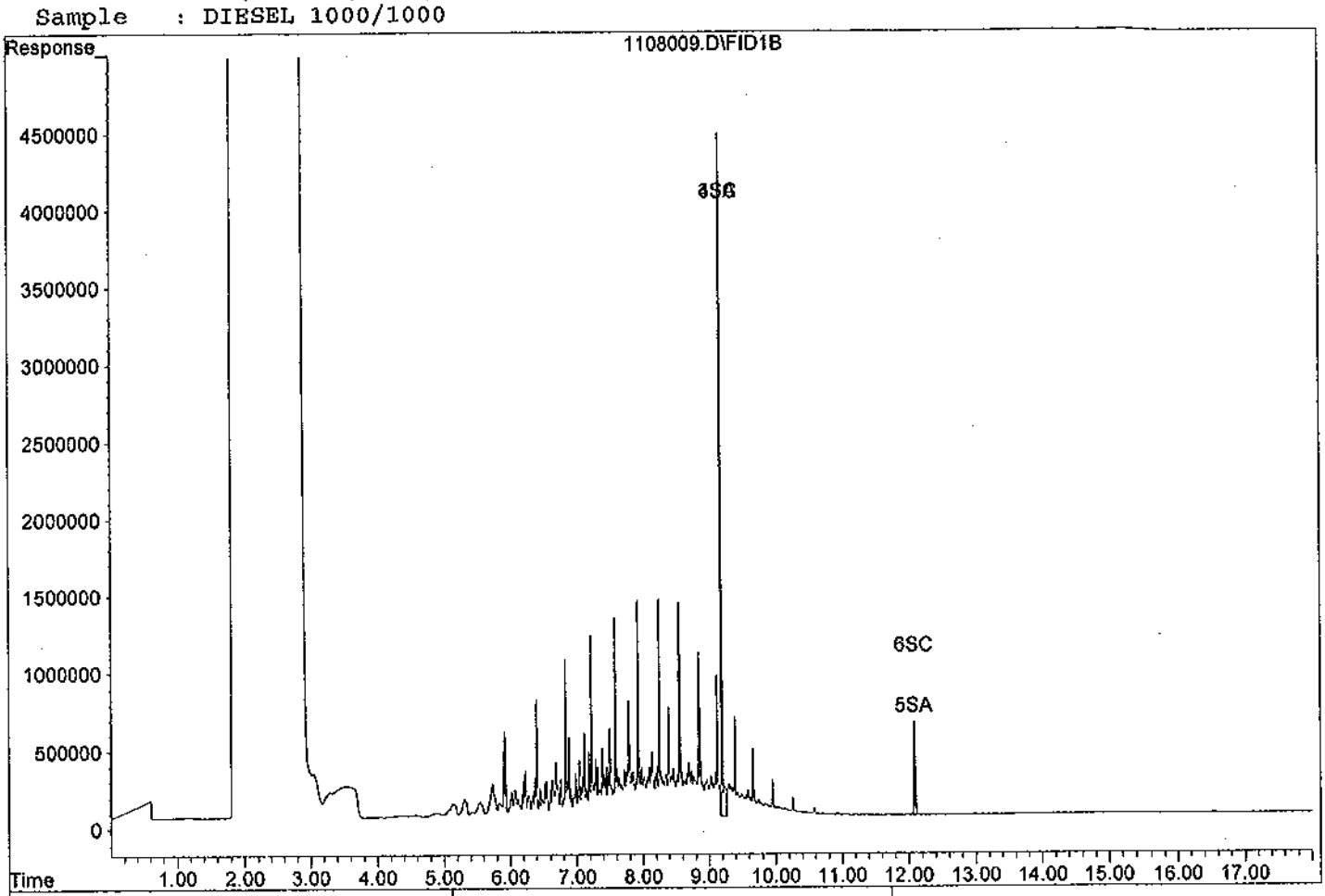
Method : G:\APOLLO\DATA\111108\TPH8S15.M (Chemstation Integrator)
Title : Diesel
Last Update : Wed Nov 30 11:52:46 2011
Response via : Multiple Level Calibration

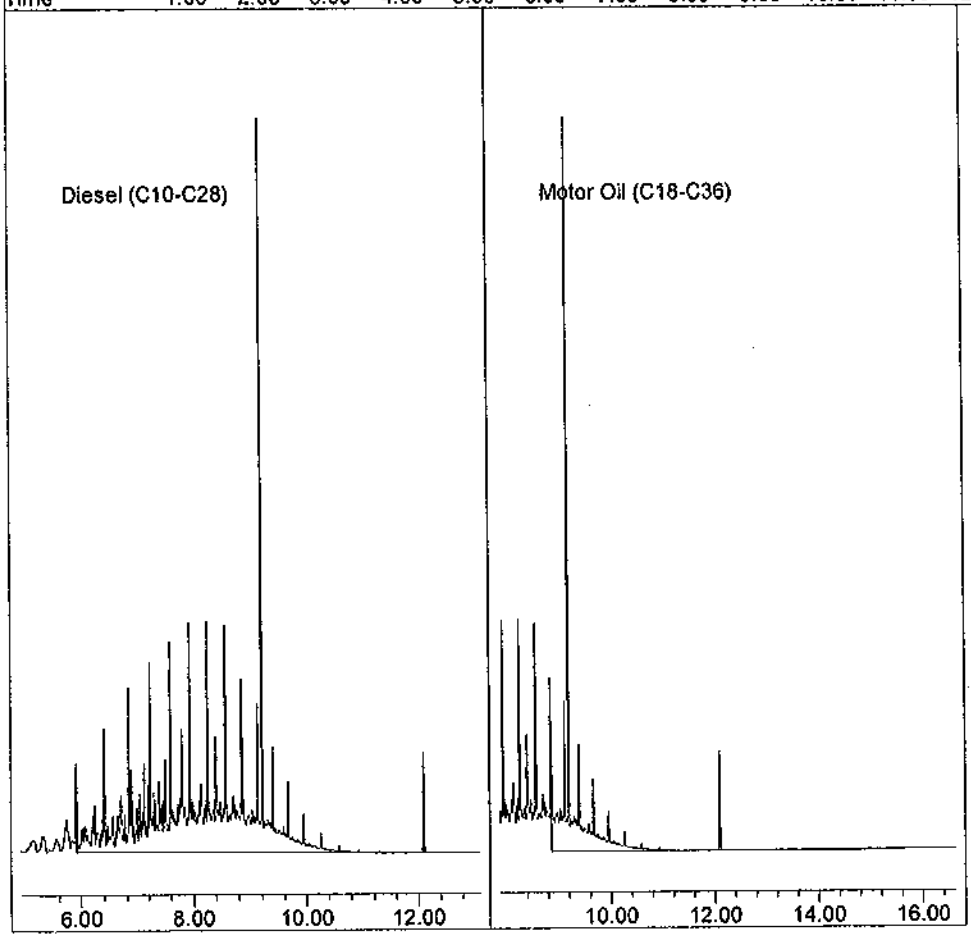
Volume Inj. : 2UL
Signal Phase : DB-5
Signal Info : FID02A

Compound	R.T.	Response	Conc	Units
System Monitoring Compounds				
3) SA Not Used(S)	9.20	38756601	57.099	ppb
Surrogate Spike 30.000	Recovery =		190.33%	
4) SC Ortho-Terphenyl(S)	9.20	38756601	60.745	ppb
Surrogate Spike 30.000	Recovery =		202.48%	
5) SA Not Used2(S)	12.09	7987688	50.659	ppb
Surrogate Spike 30.000	Recovery =		168.86%	
6) SC Octacosane(S)	12.09	7987688	34.210	ppb
Surrogate Spike 30.000	Recovery =		114.03%	
Target Compounds				
1) HATM Diesel (C10-C28)	9.01	490402243	1004.928	ppb
2) HBTM Motor Oil (C18-C36)	12.24	123592393	532.253	ppb

(f)=RT Delta > 1/2 Window 93 (m)=manual int.

Data File: G:\APOLLO\DATA\111108\1108009.D
Sample : DIESEL 1000/1000





Quantitation Report (Not Reviewed)

Data File : G:\APOLLO\DATA\111108\1108011.D Vial: 11
Acq On : 11-8-11 18:12:45 Operator: LAC
Sample : MOTOR OIL 50/1000 11/8/11 Inst : Apollo
Misc : Mix(B) Multiplr: 1.00
IntFile : events.e
Quant Time: Nov 16 9:53 2011 Quant Results File: TPH8S15.RES

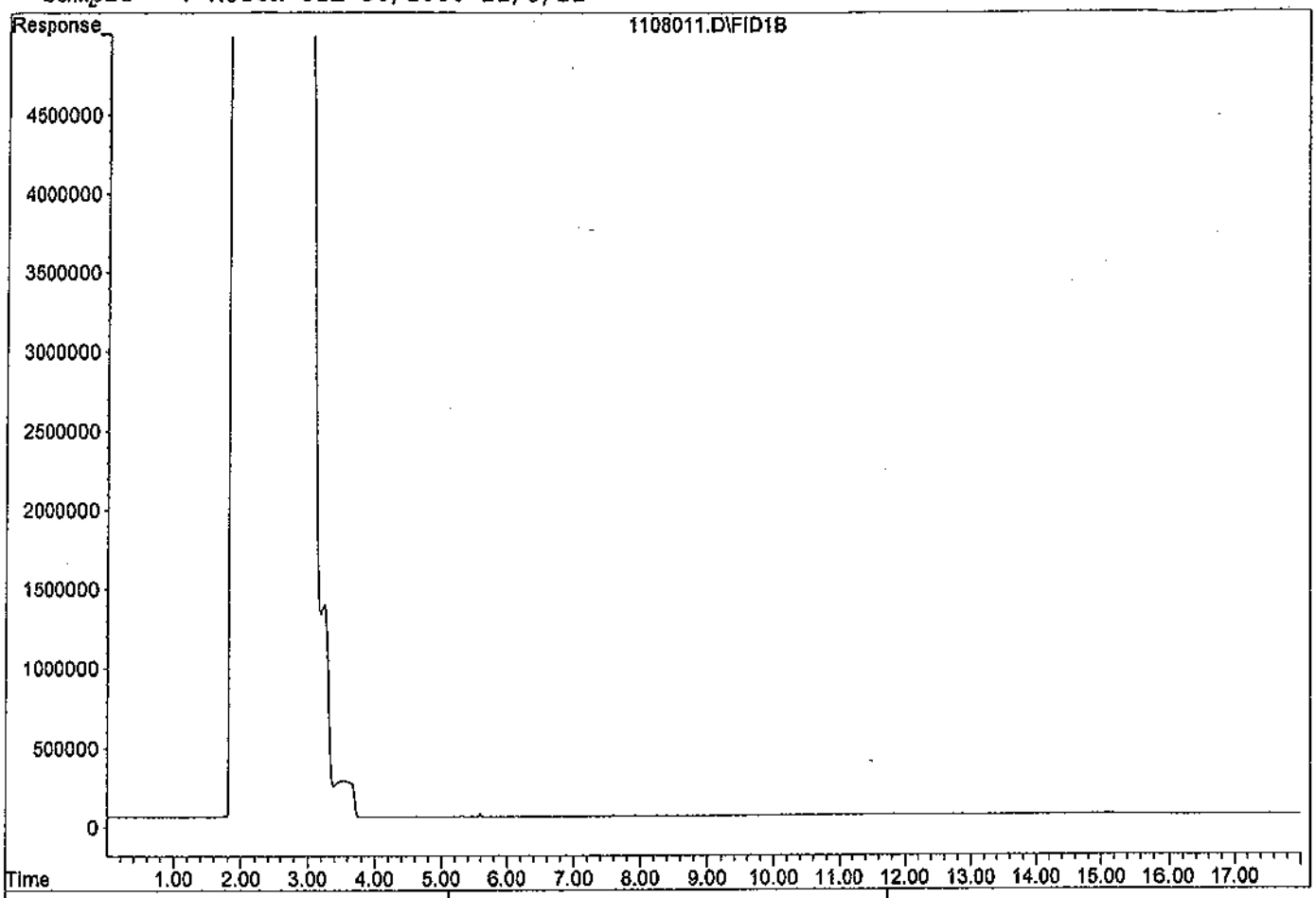
Method : G:\APOLLO\DATA\111115\TPH8S15.M (Chemstation Integrator)
Title : Diesel
Last Update : Thu Nov 17 09:41:49 2011
Response via : Multiple Level Calibration

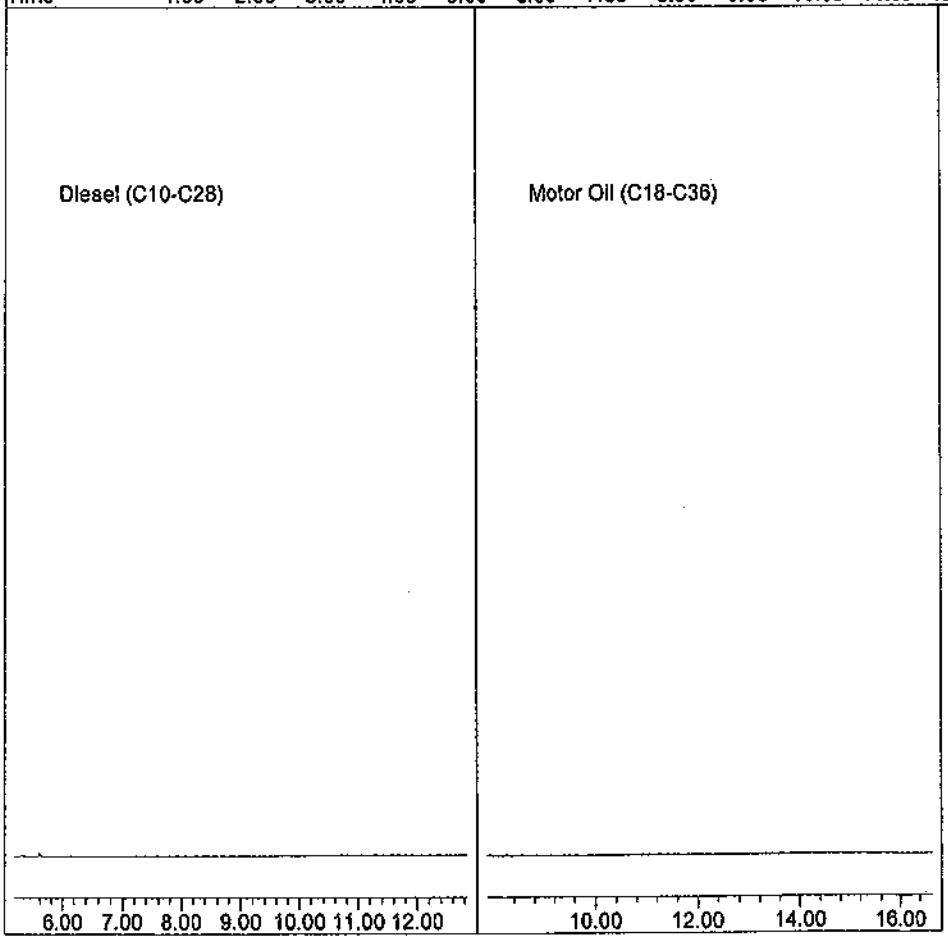
Volume Inj. : 2UL
Signal Phase : DB-5
Signal Info : FID02A

Compound	R.T.	Response	Conc Units
System Monitoring Compounds			
Target Compounds			
2) HBTM Motor Oil (C18-C36)	12.24	14043686	169.078 ppb

Quantitation Report

Data File: G:\APOLLO\DATA\111108\1108011.D
Sample : MOTOR OIL 50/1000 11/8/11





Quantitation Report (Not Reviewed)

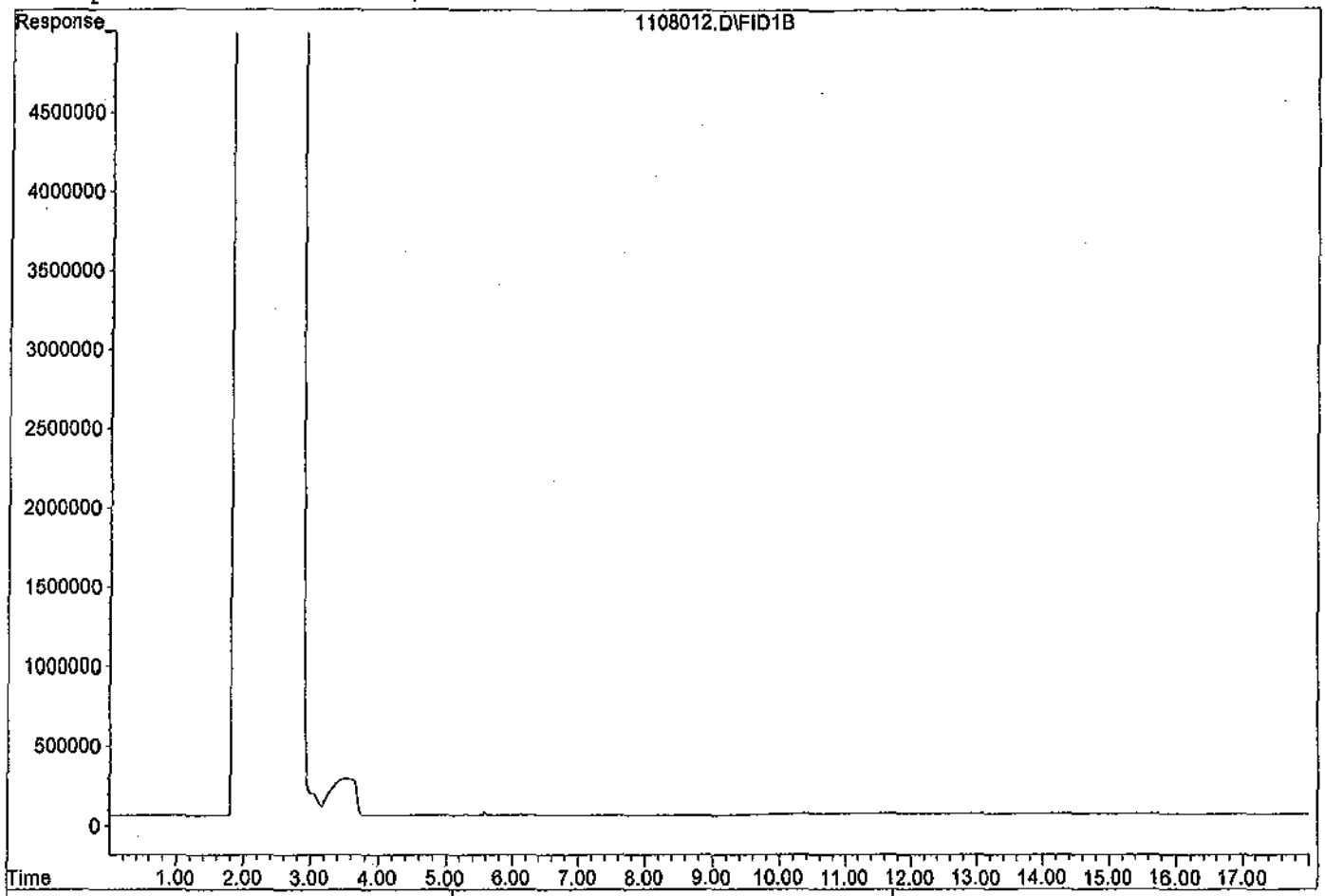
Data File : G:\APOLLO\DATA\111108\1108012.D Vial: 12
Acq On : 11-8-11 18:36:14 Operator: LAC
Sample : MOTOR OIL 100/1000 Inst : Apollo
Misc : Mix(B) Multiplr: 1.00
IntFile : events.e
Quant Time: Nov 16 9:53 2011 Quant Results File: TPH8S15.RES

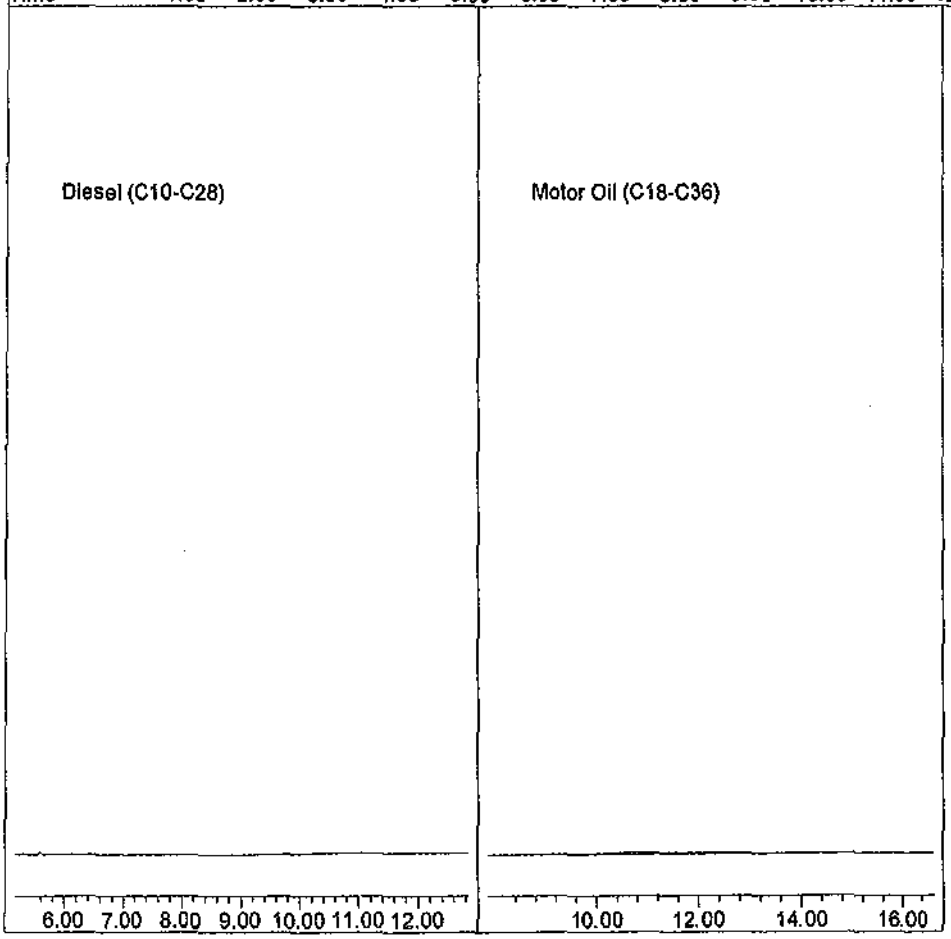
Method : G:\APOLLO\DATA\111115\TPH8S15.M (Chemstation Integrator)
Title : Diesel
Last Update : Thu Nov 17 09:41:49 2011
Response via : Multiple Level Calibration

Volume Inj. : 2UL
Signal Phase : DB-5
Signal Info : FID02A

Compound	R.T.	Response	Conc Units
System Monitoring Compounds			
Target Compounds			
2) HBTM Motor Oil (C18-C36)	12.24	19926419	239.903 ppb

Data File: G:\APOLLO\DATA\111108\1108012.D
Sample : MOTOR OIL 100/1000





Quantitation Report (Not Reviewed)

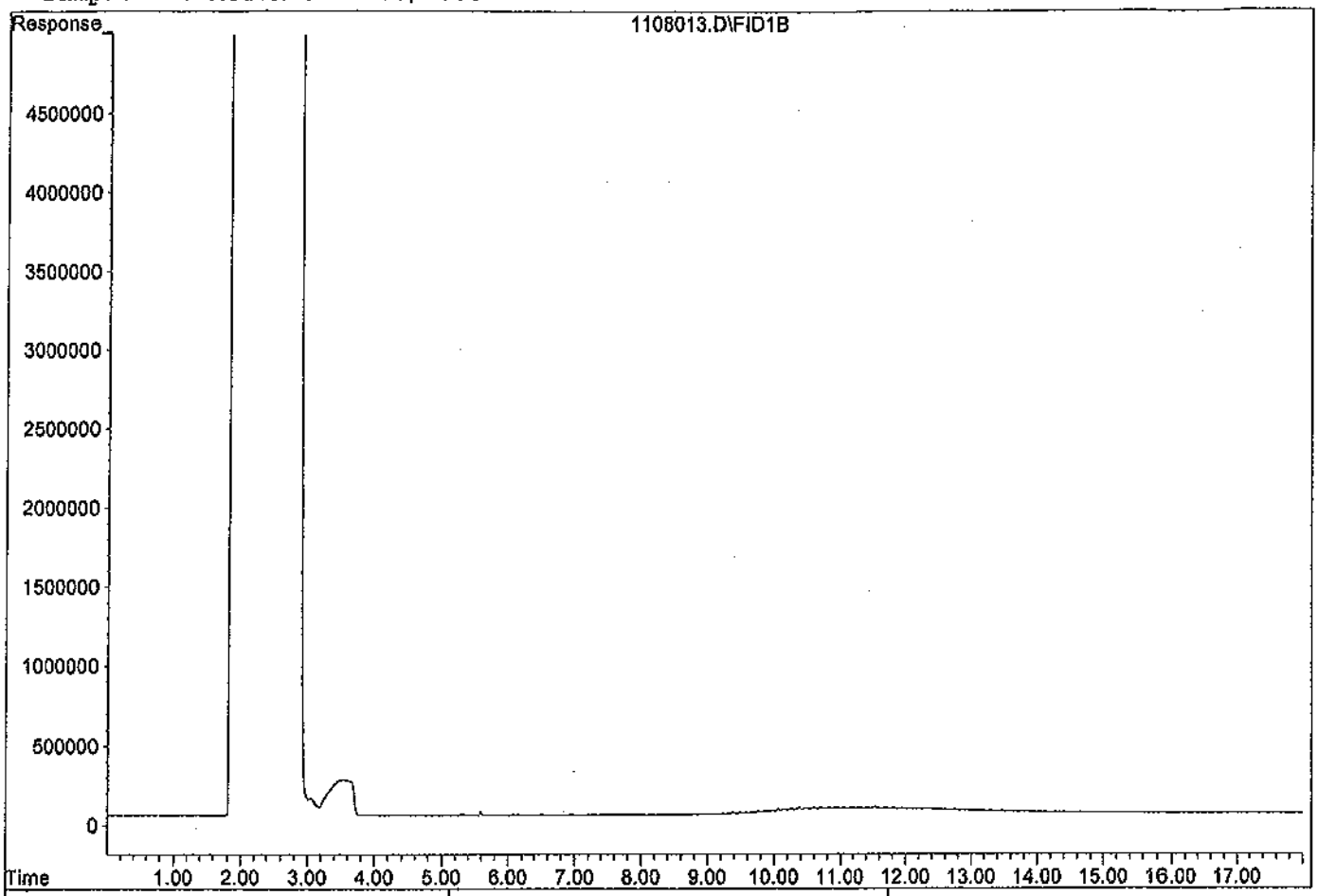
Data File : G:\APOLLO\DATA\111108\1108013.D Vial: 13
Acq On : 11-8-11 18:59:47 Operator: LAC
Sample : MOTOR OIL 400/1000 Inst : Apollo
Misc : Mix(B) Multiplr: 1.00
IntFile : events.e
Quant Time: Nov 16 9:53 2011 Quant Results File: TPH8S15.RES

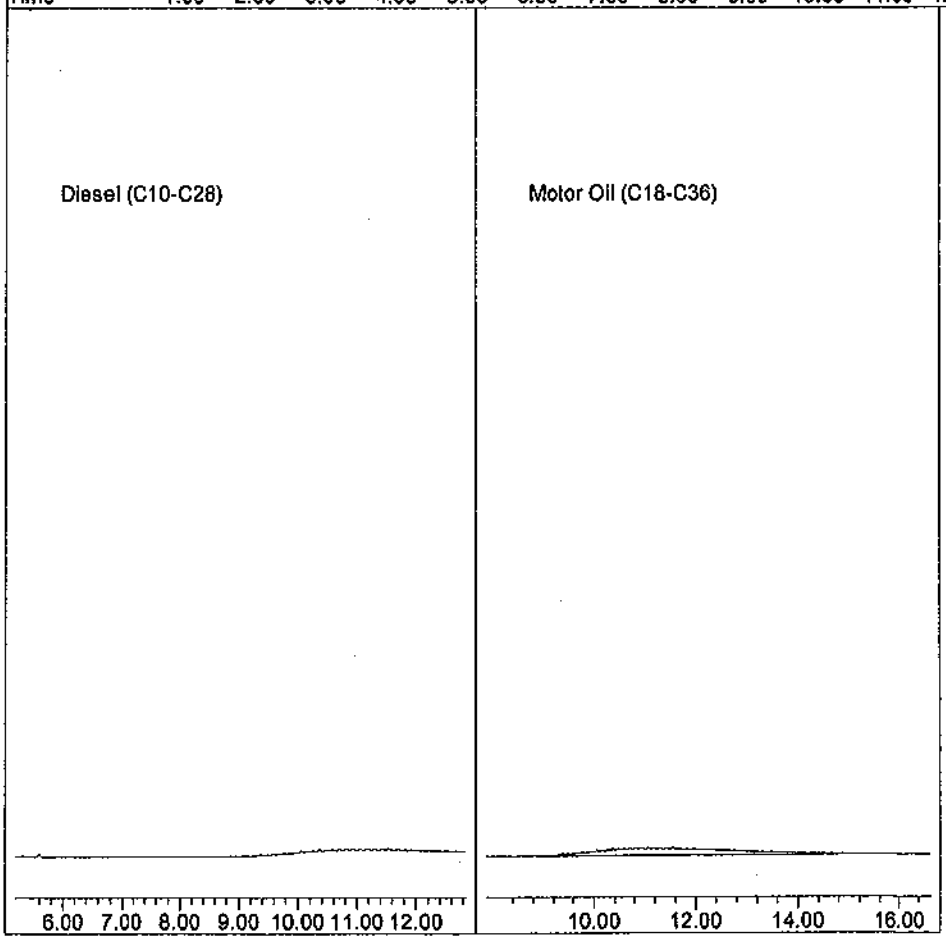
Method : G:\APOLLO\DATA\111115\TPH8S15.M (Chemstation Integrator)
Title : Diesel
Last Update : Thu Nov 17 09:41:49 2011
Response via : Multiple Level Calibration

Volume Inj. : 2UL
Signal Phase : DB-5
Signal Info : FID02A

Compound	R.T.	Response	Conc	Units
System Monitoring Compounds				
Target Compounds				
2) HBTM Motor Oil (C18-C36)	12.24	83351892	1003.512	ppb

Data File: G:\APOLLO\DATA\111108\1108013.D
Sample : MOTOR OIL 400/1000





Quantitation Report (Not Reviewed)

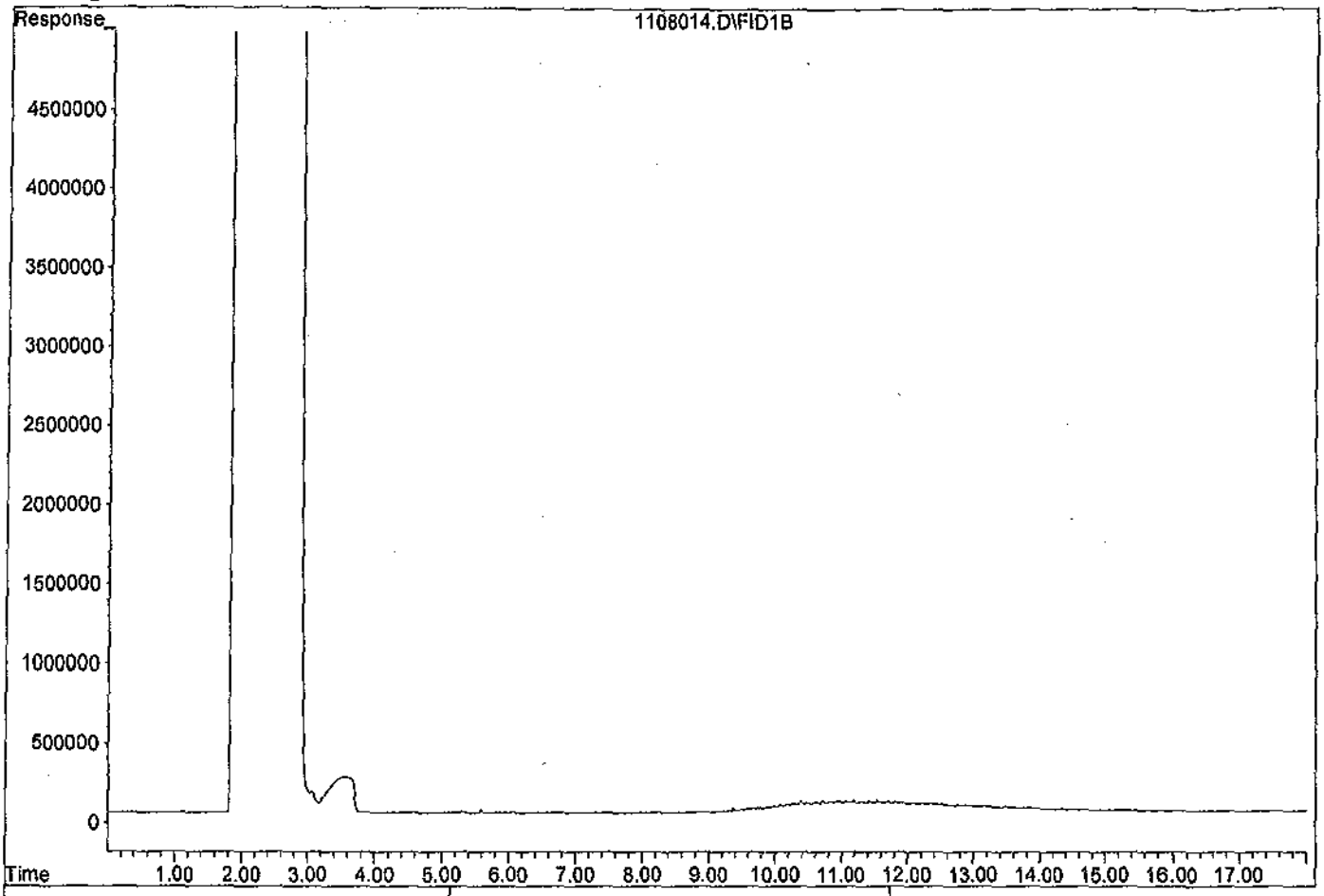
Data File : G:\APOLLO\DATA\111108\1108014.D Vial: 14
Acq On : 11-8-11 19:23:20 Operator: LAC
Sample : MOTOR OIL 600/1000 Inst : Apollo
Misc : Mix(B) Multiplr: 1.00
IntFile : events.e
Quant Time: Nov 16 9:54 2011 Quant Results File: TPH8S15.RES

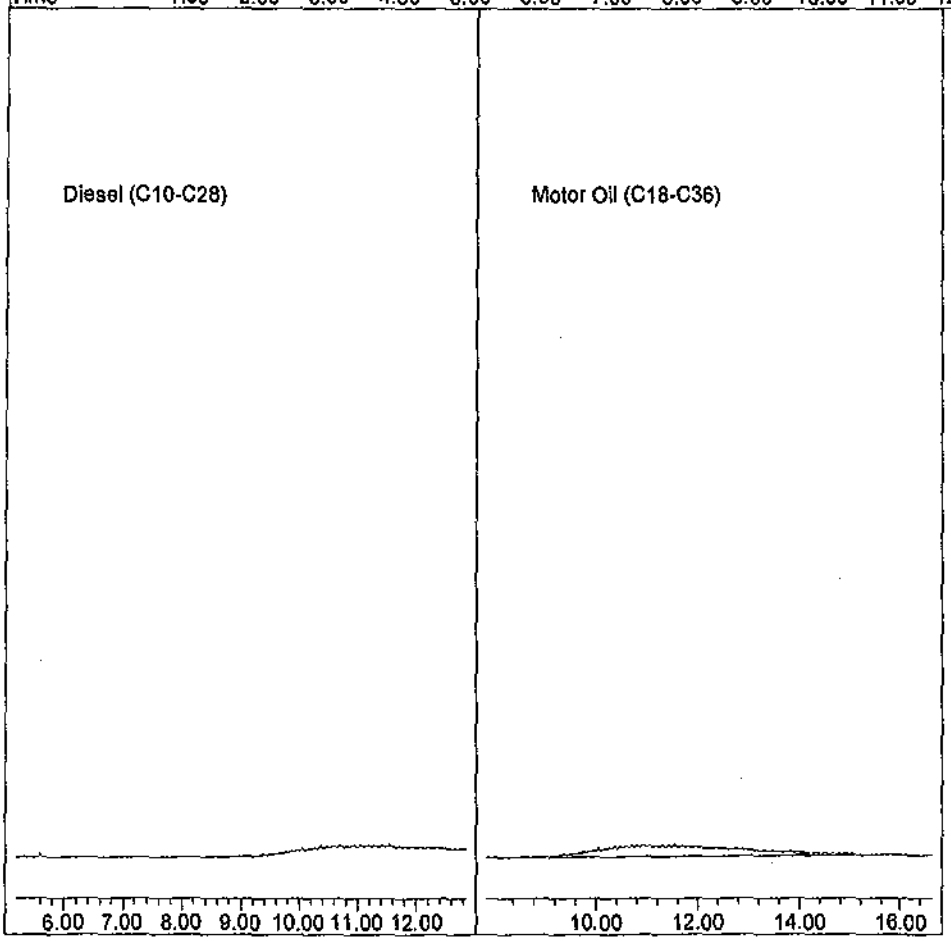
Method : G:\APOLLO\DATA\111115\TPH8S15.M (Chemstation Integrator)
Title : Diesel
Last Update : Thu Nov 17 09:41:49 2011
Response via : Multiple Level Calibration

Volume Inj. : 2UL
Signal Phase : DB-5
Signal Info : FID02A

Compound	R.T.	Response	Conc	Units
System Monitoring Compounds				
Target Compounds				
2) HBTM Motor Oil (C18-C36)	12.24	133423372	1606.346	ppb

Data File: G:\APOLLO\DATA\111108\1108014.D
Sample : MOTOR OIL 600/1000





Quantitation Report (Not Reviewed)

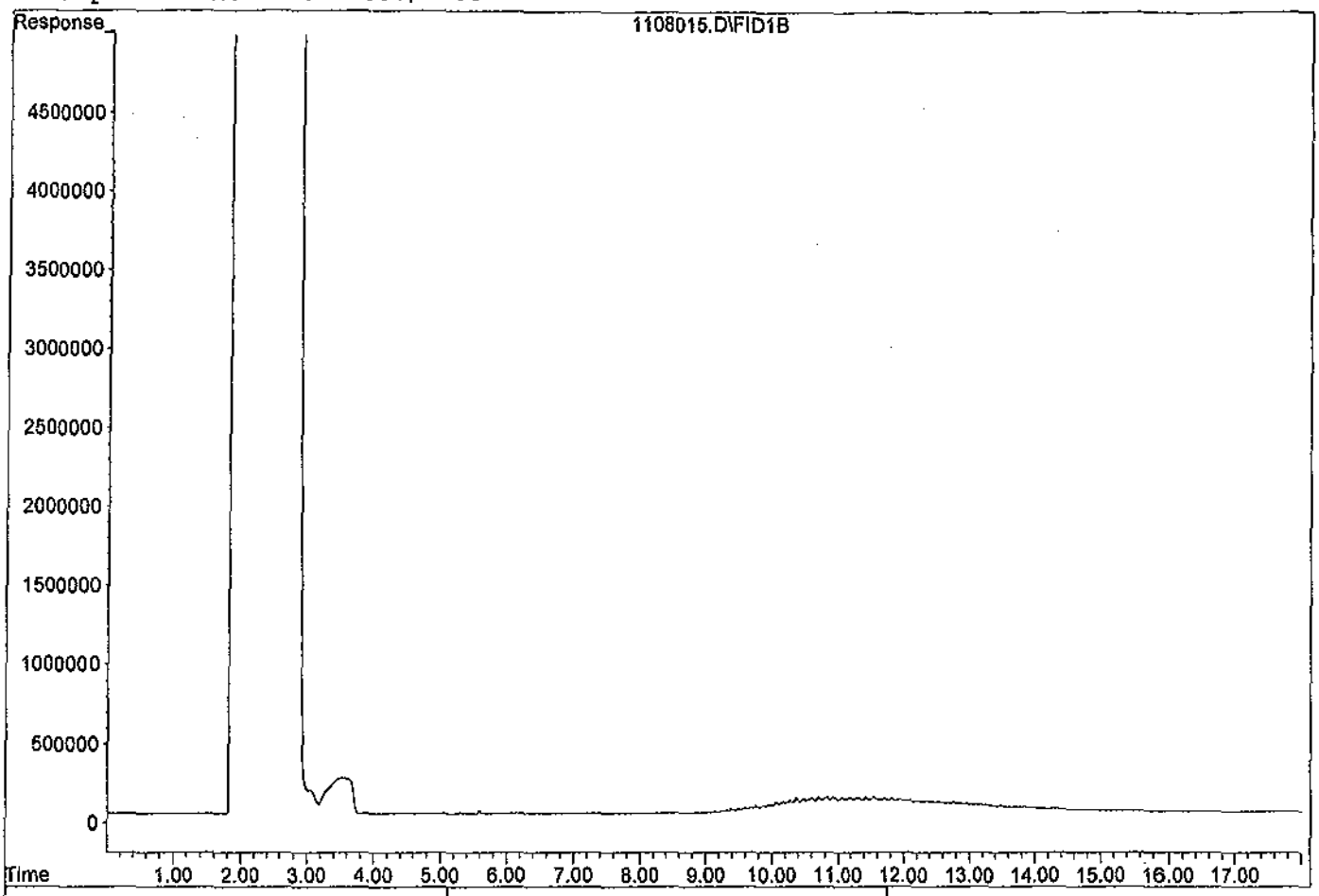
Data File : G:\APOLLO\DATA\111108\1108015.D Vial: 15
Acq On : 11-8-11 19:46:53 Operator: LAC
Sample : MOTOR OIL 800/1000 Inst : Apollo
Misc : Mix(B) Multiplr: 1.00
IntFile : events.e
Quant Time: Nov 16 9:54 2011 Quant Results File: TPH8S15.RES

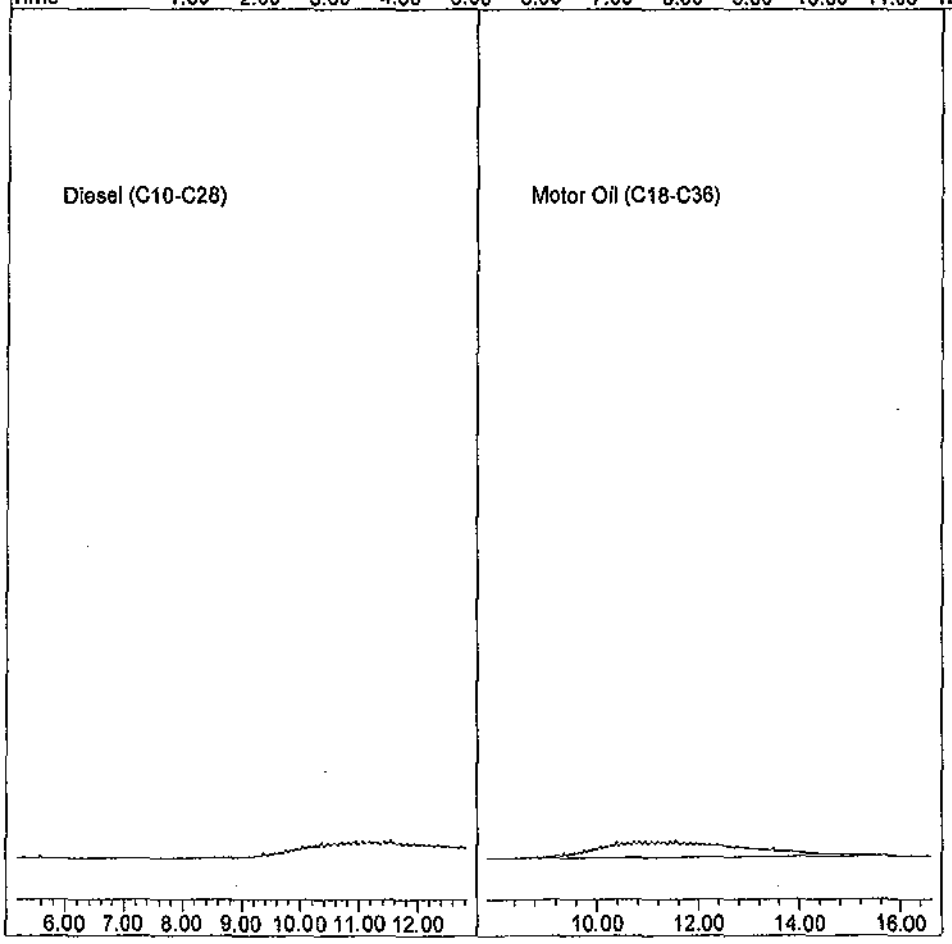
Method : G:\APOLLO\DATA\111115\TPH8S15.M (Chemstation Integrator)
Title : Diesel
Last Update : Thu Nov 17 09:41:49 2011
Response via : Multiple Level Calibration

Volume Inj. : 2UL
Signal Phase : DB-5
Signal Info : FID02A

Compound	R.T.	Response	Conc	Units
System Monitoring Compounds				
Target Compounds				
2) HBTM Motor Oil (C18-C36)	12.24	185280557	2230.679	ppb

Data File: G:\APOLLO\DATA\111108\1108015.D
Sample : MOTOR OIL 800/1000





Quantitation Report (Not Reviewed)

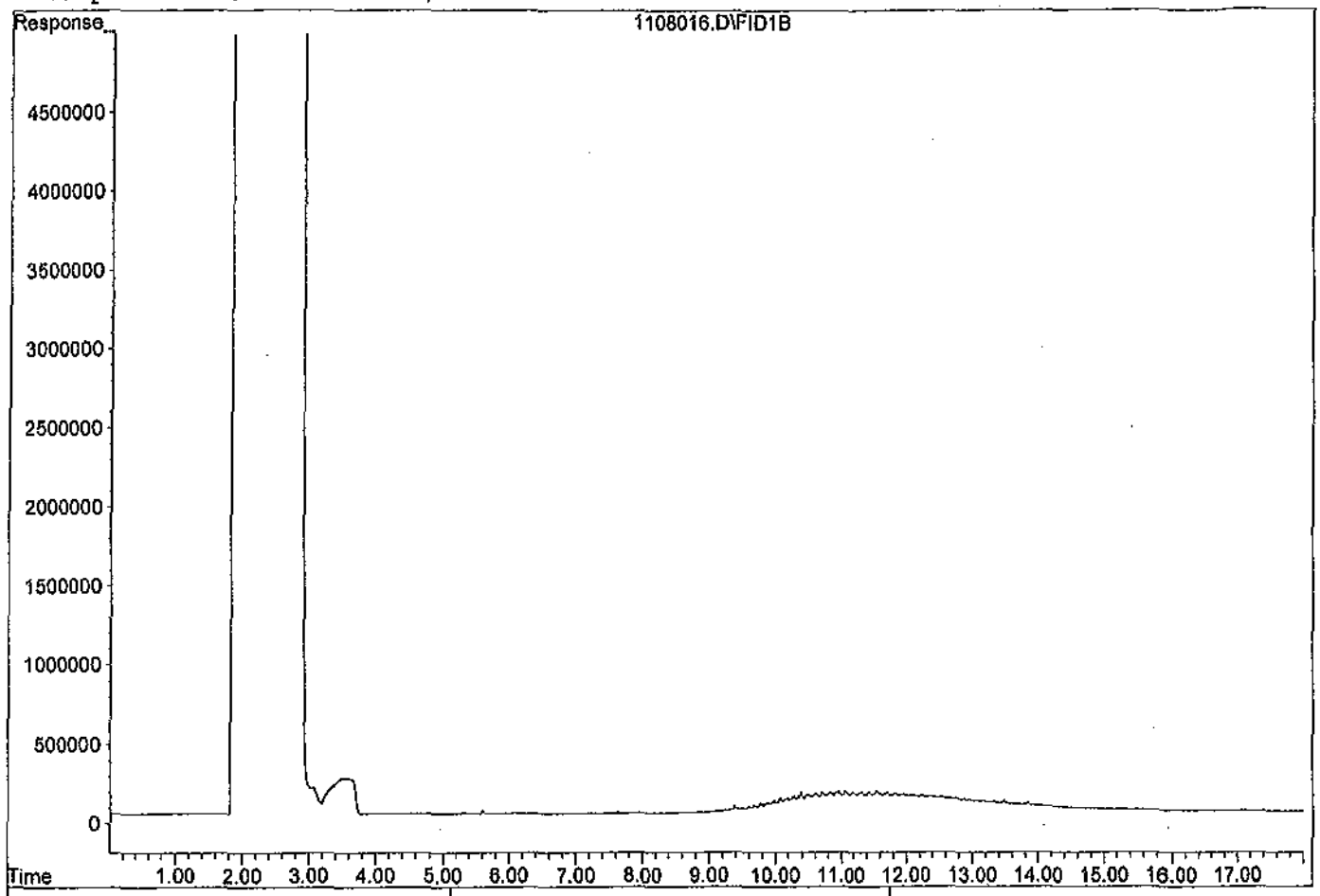
Data File : G:\APOLLO\DATA\111108\1108016.D Vial: 16
Acq On : 11-8-11 20:10:21 Operator: LAC
Sample : MOTOR OIL 1000/1000 Inst : Apollo
Misc : Mix(B) Multiplr: 1.00
IntFile : events.e
Quant Time: Nov 16 9:54 2011 Quant Results File: TPH8S15.RES

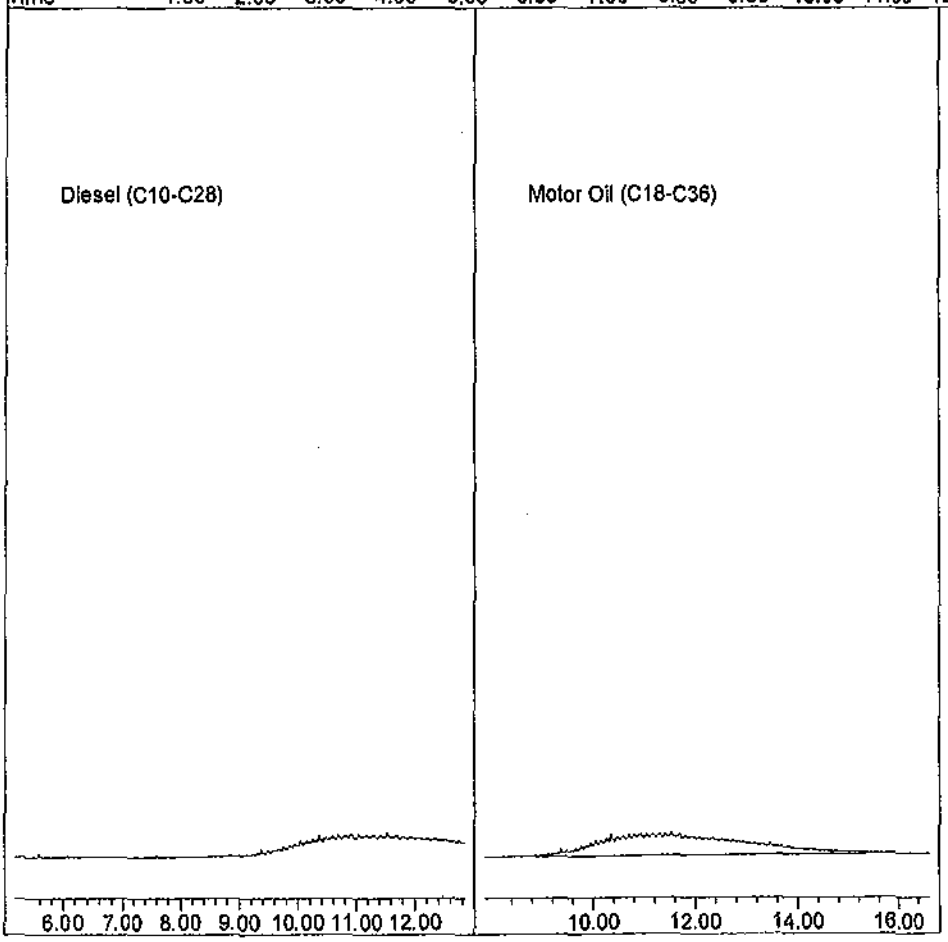
Method : G:\APOLLO\DATA\111115\TPH8S15.M (Chemstation Integrator)
Title : Diesel
Last Update : Thu Nov 17 09:41:49 2011
Response via : Multiple Level Calibration

Volume Inj. : 2UL
Signal Phase : DB-5
Signal Info : FID02A

Compound	R.T.	Response	Conc	Units
System Monitoring Compounds				
Target Compounds				
2) HBTM Motor Oil (C18-C36)	12.24	250746792	3018.857	ppb

Data File: G:\APOLLO\DATA\111108\1108016.D
Sample : MOTOR OIL 1000/1000





Quantitation Report (Not Reviewed)

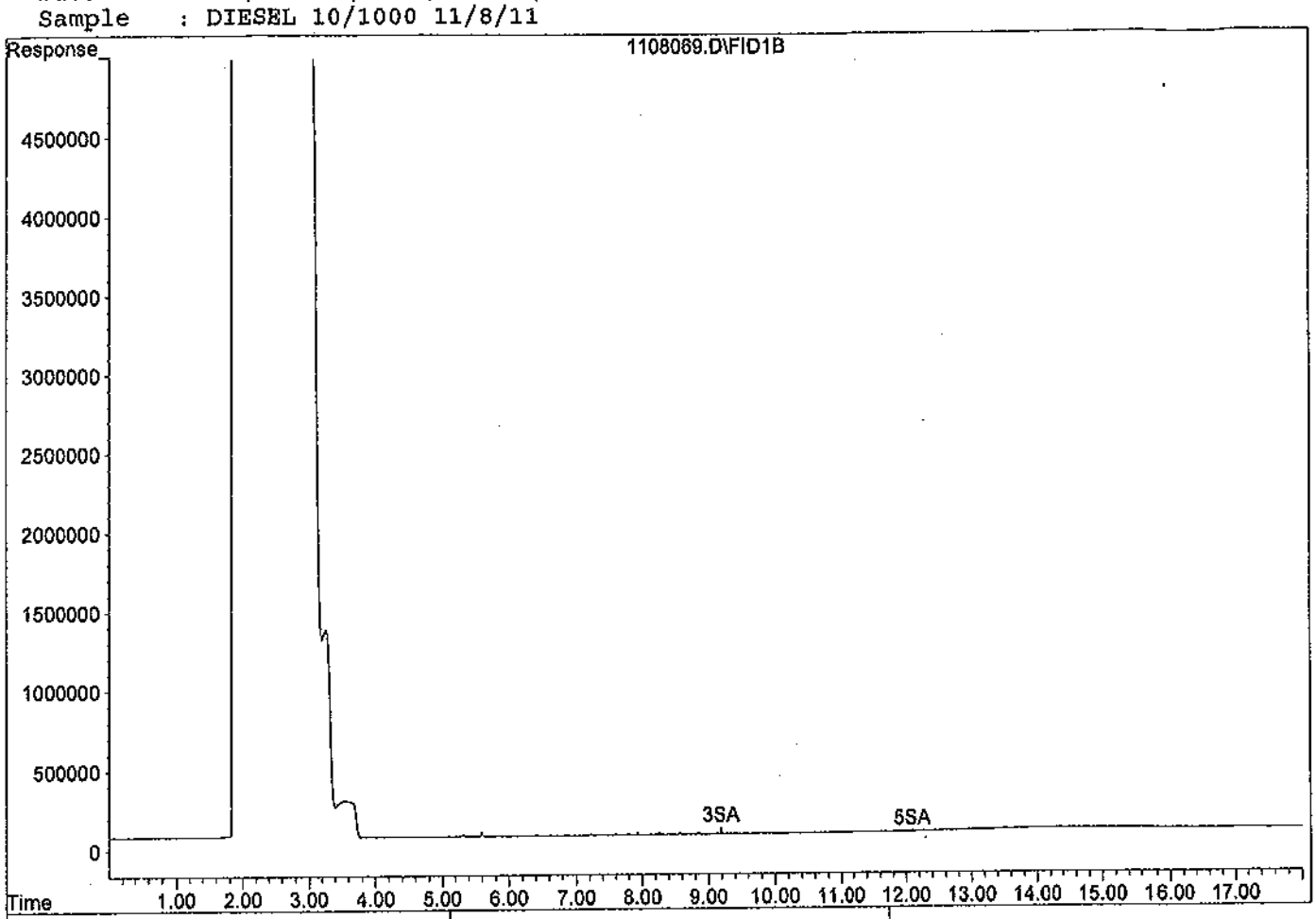
Data File : G:\APOLLO\DATA\111108\1108069.D Vial: 69
Acq On : 11-9-11 17:18:58 Operator: LAC
Sample : DIESEL 10/1000 11/8/11 Inst : Apollo
Misc : Mix(A) Multiplr: 1.00
IntFile : events.e
Quant Time: Nov 16 9:53 2011 Quant Results File: TPH8S15.RES

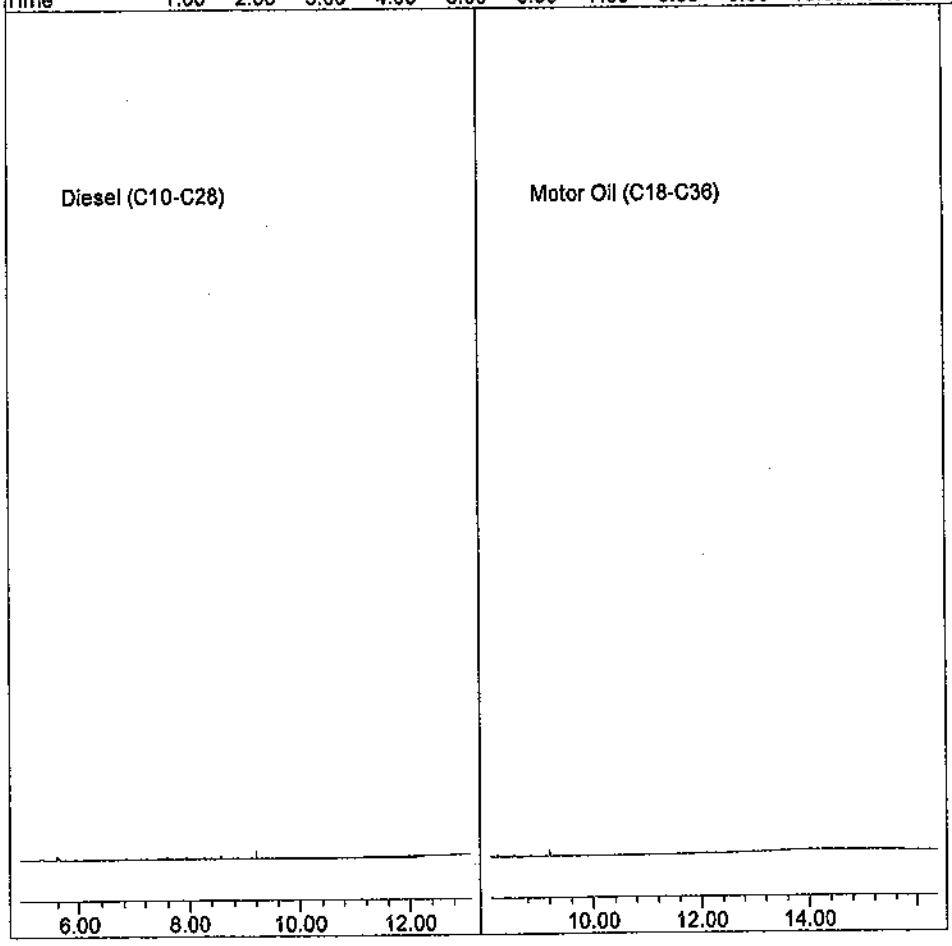
Method : G:\APOLLO\DATA\111115\TPH8S15.M (Chemstation Integrator)
Title : Diesel
Last Update : Thu Nov 17 09:41:49 2011
Response via : Multiple Level Calibration

Volume Inj. : 2UL
Signal Phase : DB-5
Signal Info : FID02A

Compound	R.T.	Response	Conc Units
System Monitoring Compounds			
3) SA Not Used(S)	9.20	302444	0.297 ppb
Surrogate Spike 30.000	Recovery	=	0.99%
5) SA Not Used2(S)	12.10	625179	2.122 ppb
Surrogate Spike 30.000	Recovery	=	7.07%
Target Compounds			
1) HATM Diesel (C10-C28)	9.01	12262633	1055.198 ppb

Data File: G:\APOLLO\DATA\111108\1108069.D
Sample : DIESEL 10/1000 11/8/11





Quantitation Report (Not Reviewed)

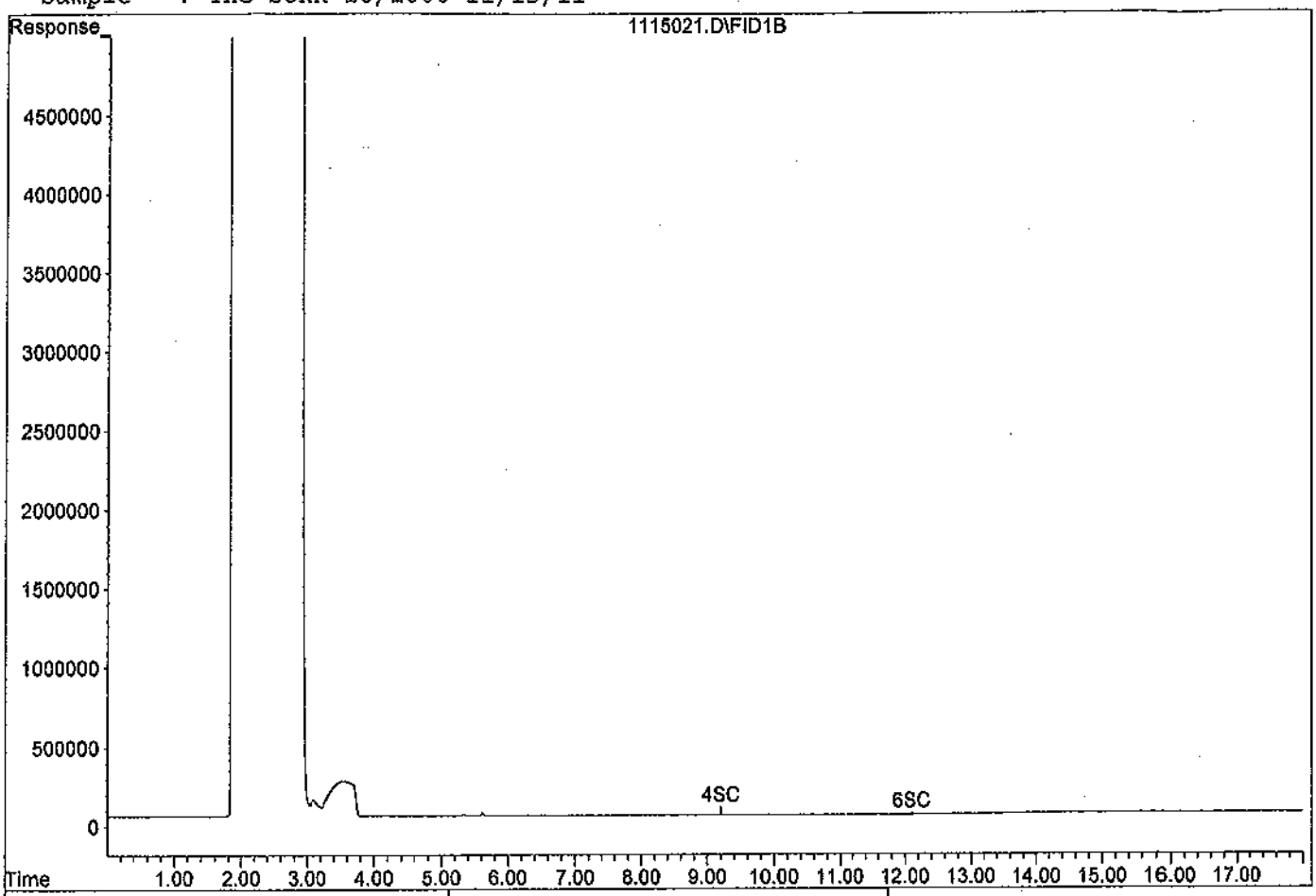
Data File : G:\APOLLO\DATA\111115\1115021.D Vial: 21
Acq On : 11-15-11 18:21:35 Operator: LAC
Sample : THC SURR 10/1000 11/15/11 Inst : Apollo
Misc : Mix(C) Multiplr: 1.00
IntFile : events.e
Quant Time: Nov 16 14:58 2011 Quant Results File: TPH8S15.RES

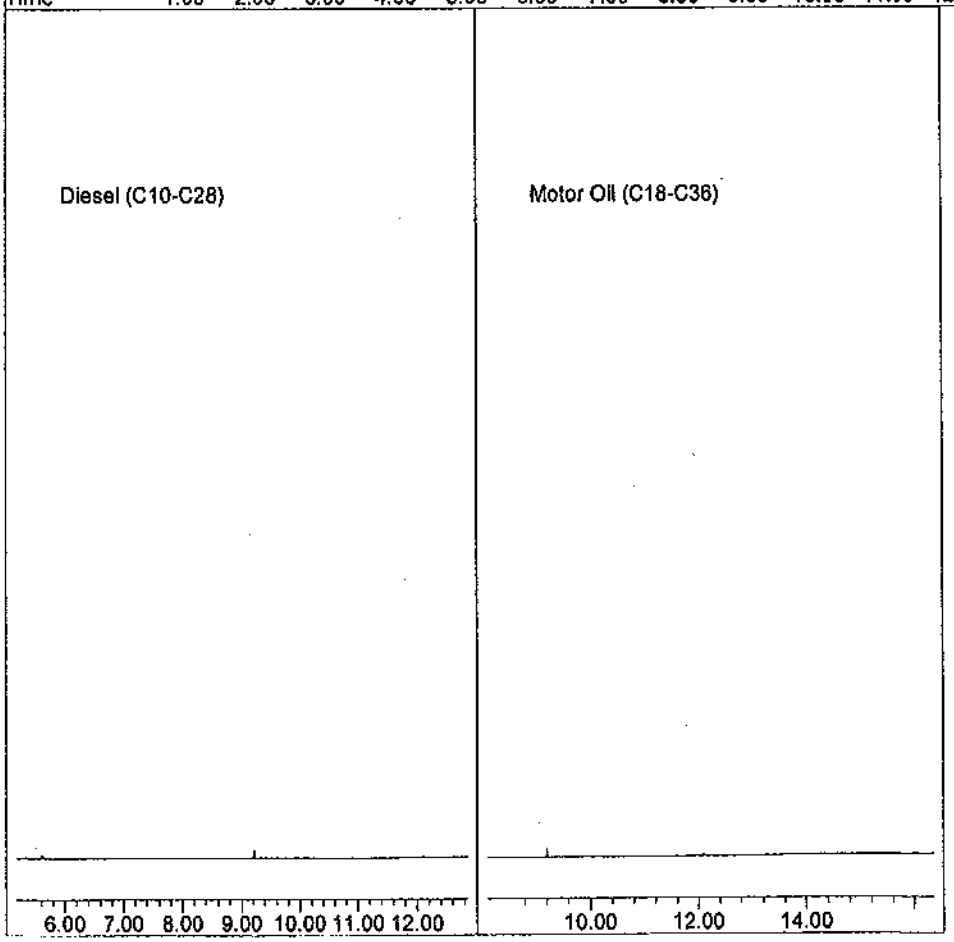
Method : G:\APOLLO\DATA\111115\TPH8S15.M (Chemstation Integrator)
Title : Diesel
Last Update : Thu Nov 17 09:41:49 2011
Response via : Multiple Level Calibration

Volume Inj. : 2UL
Signal Phase : DB-5
Signal Info : FID02A

Compound	R.T.	Response	Conc Units
System Monitoring Compounds			
4) SC Ortho-Terphenyl(S)	9.21	356915	0.559 ppb
Surrogate Spike 30.000	Recovery	=	1.86%
6) SC Octacosane(S)	12.10	279297	1.196 ppb
Surrogate Spike 30.000	Recovery	=	3.99%
Target Compounds			

Data File: G:\APOLLO\DATA\111115\1115021.D
Sample : THC SURR 10/1000 11/15/11





Quantitation Report (Not Reviewed)

Data File : G:\APOLLO\DATA\111115\1115022.D Vial: 22
Acq On : 11-15-11 18:45:31 Operator: LAC
Sample : THC SURR 100/1000 Inst : Apollo
Misc : Mix(C) Multiplr: 1.00
IntFile : events.e
Quant Time: Nov 16 14:58 2011 Quant Results File: TPH8S15.RES

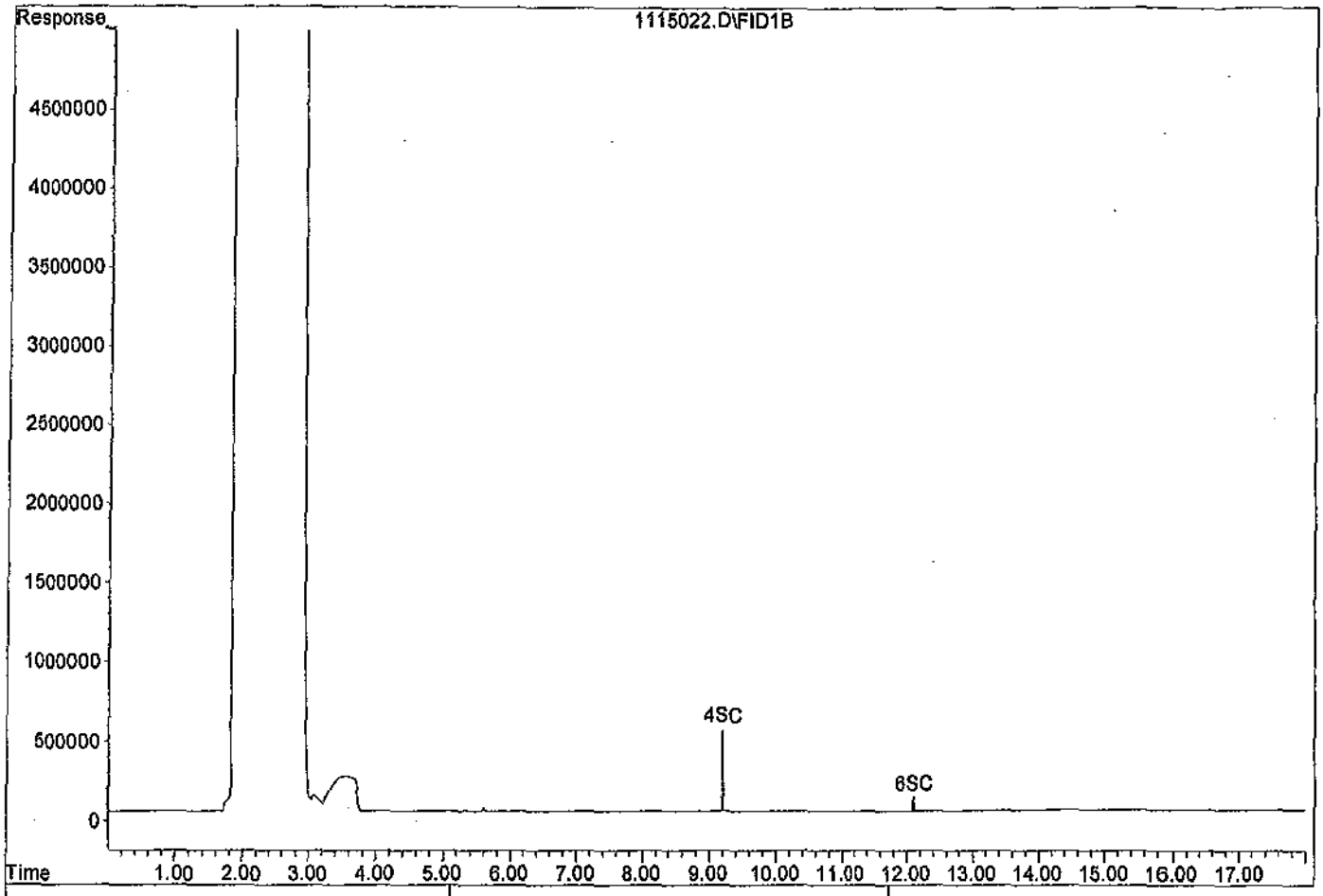
Method : G:\APOLLO\DATA\111115\TPH8S15.M (Chemstation Integrator)
Title : Diesel
Last Update : Thu Nov 17 09:41:49 2011
Response via : Multiple Level Calibration

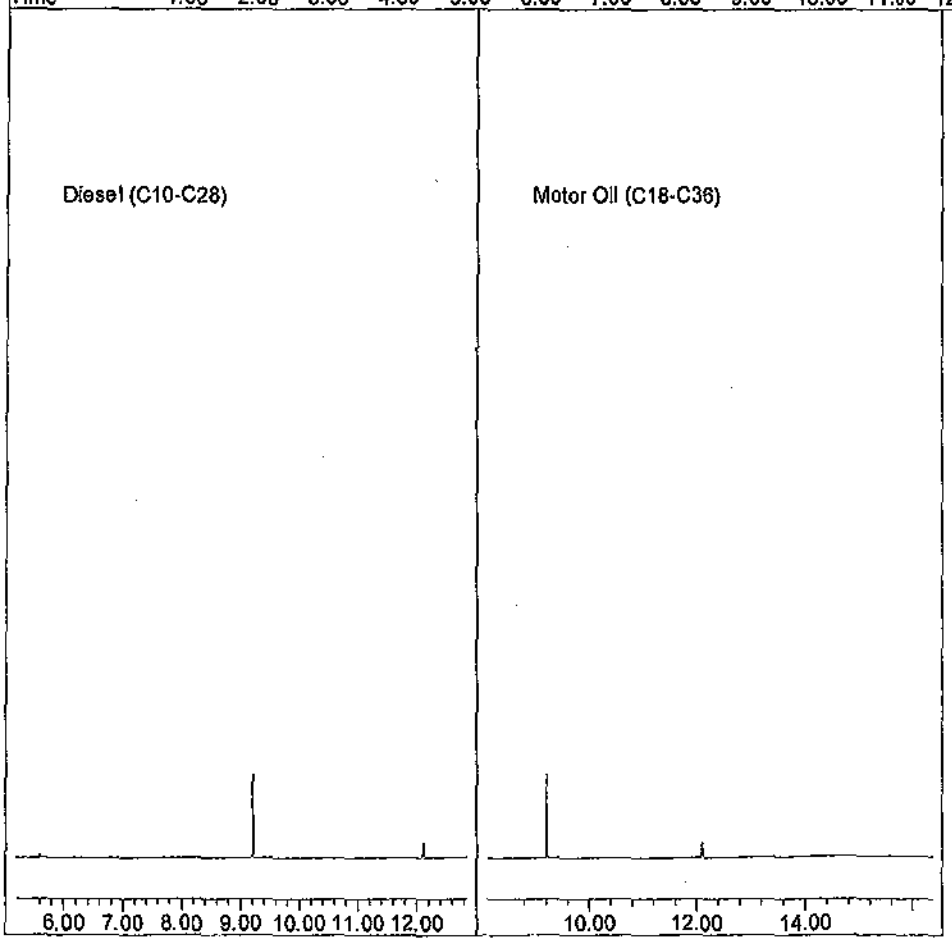
Volume Inj. : 2UL
Signal Phase : DB-5
Signal Info : FID02A

Compound	R.T.	Response	Conc	Units
System Monitoring Compounds				
4) SC Ortho-Terphenyl(S)	9.21	3207972	5.028	ppb
Surrogate Spike 30.000	Recovery =		16.76%	
6) SC Octacosane(S)	12.11	1214451	5.201	ppb
Surrogate Spike 30.000	Recovery =		17.34%	
Target Compounds				

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

Data File: G:\APOLLO\DATA\111115\1115022.D
Sample : THC SURR 100/1000





Quantitation Report (Not Reviewed)

Data File : G:\APOLLO\DATA\111115\1115023.D Vial: 23
Acq On : 11-15-11 19:09:25 Operator: LAC
Sample : THC SURR 400/1000 Inst : Apollo
Misc : Mix(C) Multiplr: 1.00
IntFile : events.e
Quant Time: Nov 16 14:58 2011 Quant Results File: TPH8S15.RES

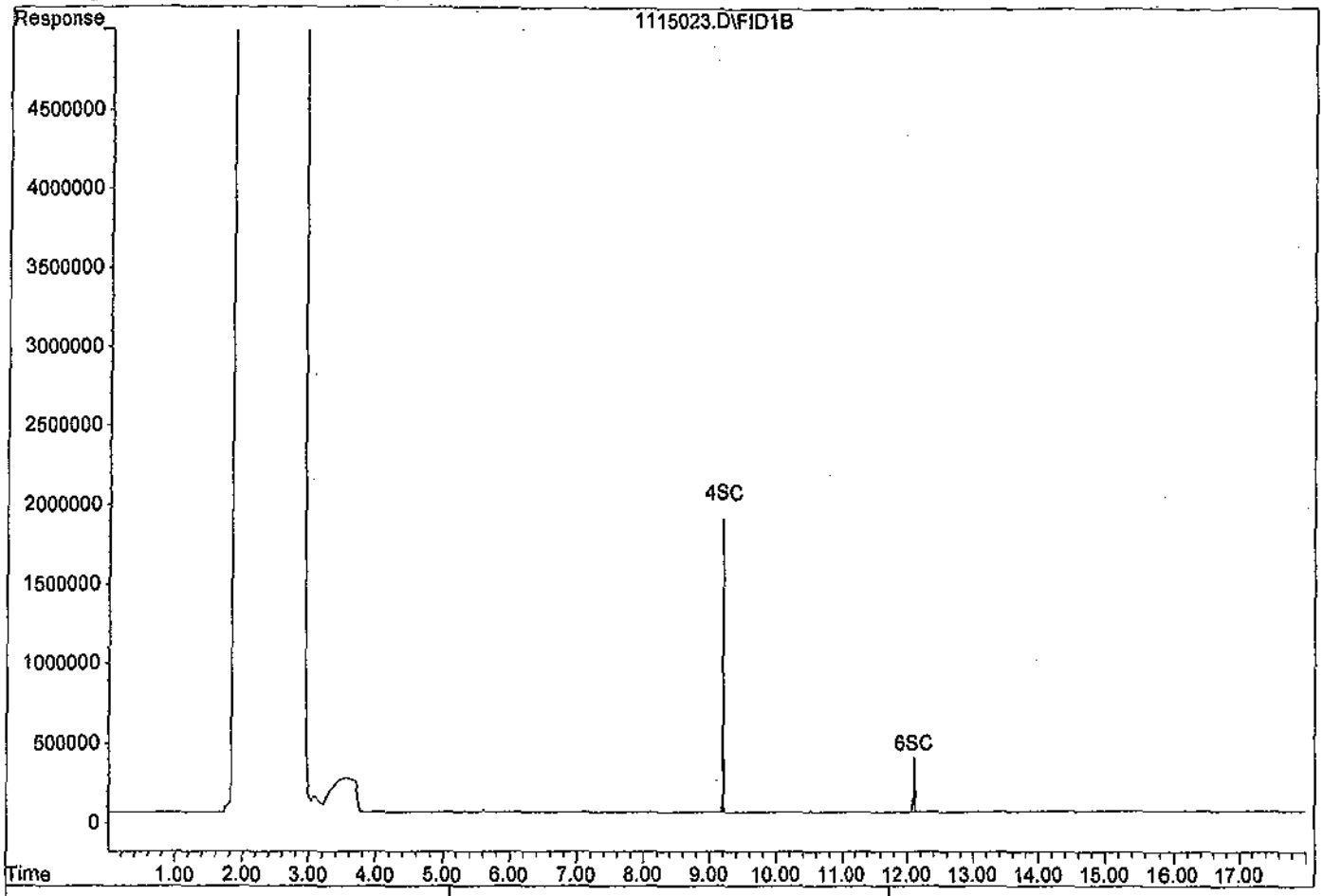
Method : G:\APOLLO\DATA\111115\TPH8S15.M (Chemstation Integrator)
Title : Diesel
Last Update : Thu Nov 17 09:41:49 2011
Response via : Multiple Level Calibration

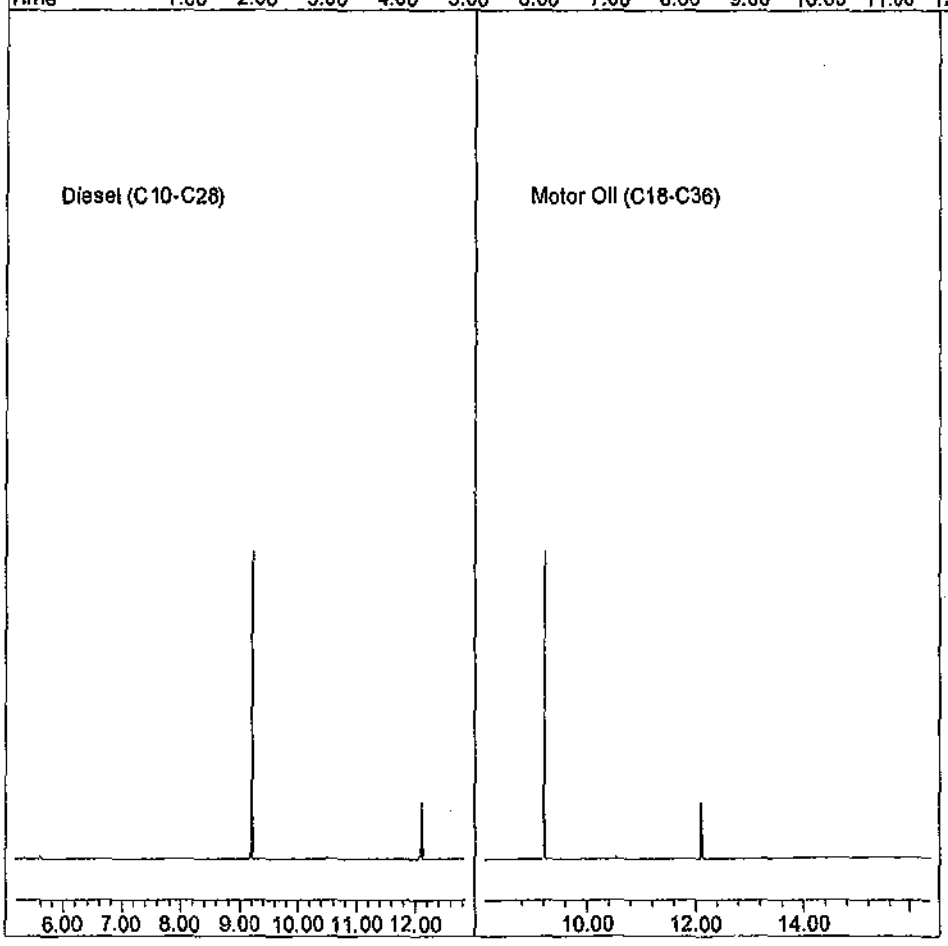
Volume Inj. : 2UL
Signal Phase : DB-5
Signal Info : FID02A

Compound	R.T.	Response	Conc	Units
System Monitoring Compounds				
4) SC Ortho-Terphenyl(S)	9.21	12023229	18.845	ppb
Surrogate Spike 30.000		Recovery =	62.82%	
6) SC Octacosane(S)	12.11	4606231	19.728	ppb
Surrogate Spike 30.000		Recovery =	65.76%	
Target Compounds				

(f)=RT Delta > 1/2 Window 113 (m)=manual int.

Data File: G:\APOLLO\DATA\111115\1115023.D
Sample : THC SURR 400/1000





Quantitation Report (Not Reviewed)

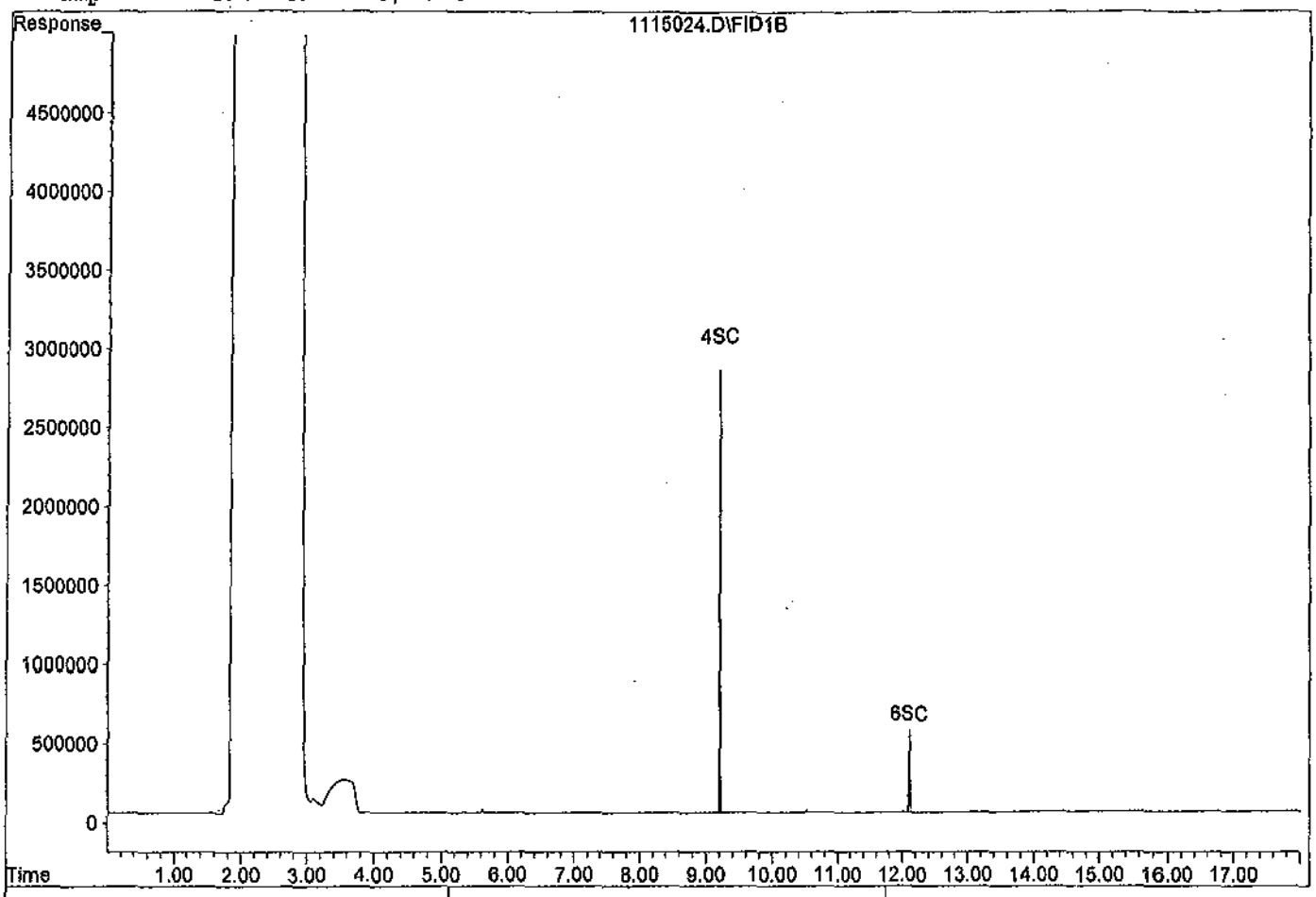
Data File : G:\APOLLO\DATA\111115\1115024.D Vial: 24
Acq On : 11-15-11 19:33:17 Operator: LAC
Sample : THC SURR 600/1000 Inst : Apollo
Misc : Mix(C) Multiplr: 1.00
IntFile : events.e
Quant Time: Nov 16 14:58 2011 Quant Results File: TPH8S15.RES

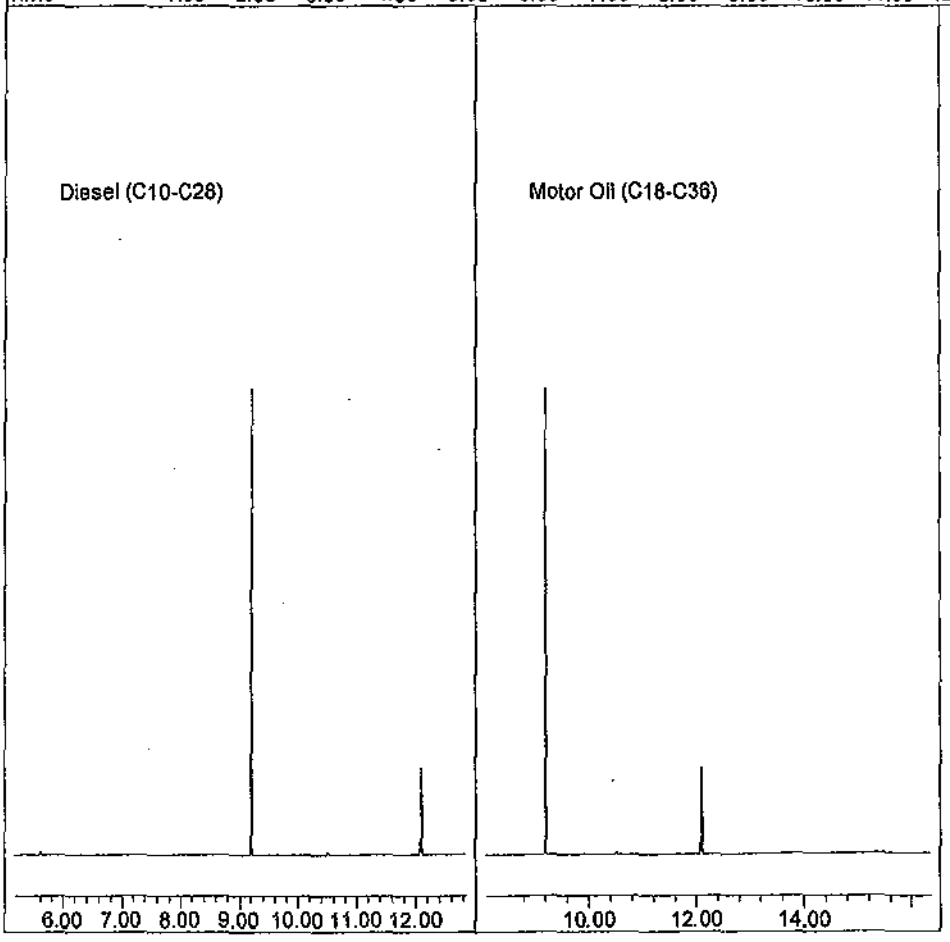
Method : G:\APOLLO\DATA\111115\TPH8S15.M (Chemstation Integrator)
Title : Diesel
Last Update : Thu Nov 17 09:41:49 2011
Response via : Multiple Level Calibration

Volume Inj. : 2UL
Signal Phase : DB-5
Signal Info : FID02A

Compound	R.T.	Response	Conc	Units
System Monitoring Compounds				
4) SC Ortho-Terphenyl(S)	9.21	18244401	28.595	ppb
Surrogate Spike 30.000	Recovery =		95.32%	
6) SC Octacosane(S)	12.11	6794679	29.101	ppb
Surrogate Spike 30.000	Recovery =		97.00%	
Target Compounds				

Data File: G:\APOLLO\DATA\111115\1115024.D
Sample : THC SURR 600/1000





Quantitation Report (Not Reviewed)

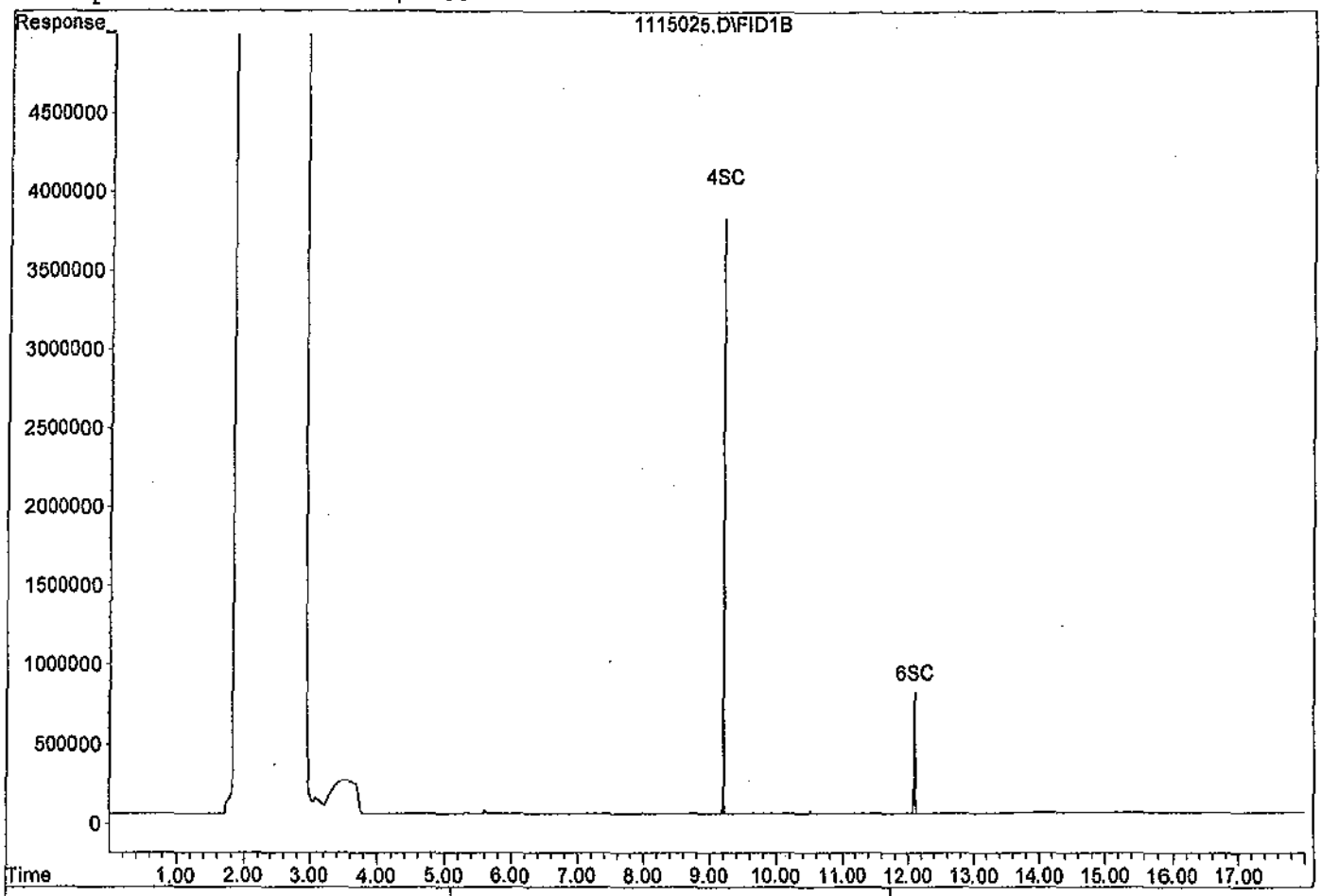
Data File : G:\APOLLO\DATA\111115\1115025.D Vial: 25
Acq On : 11-15-11 19:57:06 Operator: LAC
Sample : THC SURR 800/1000 Inst : Apollo
Misc : Mix(C) Multiplr: 1.00
IntFile : events.e
Quant Time: Nov 16 14:58 2011 Quant Results File: TPH8S15.RES

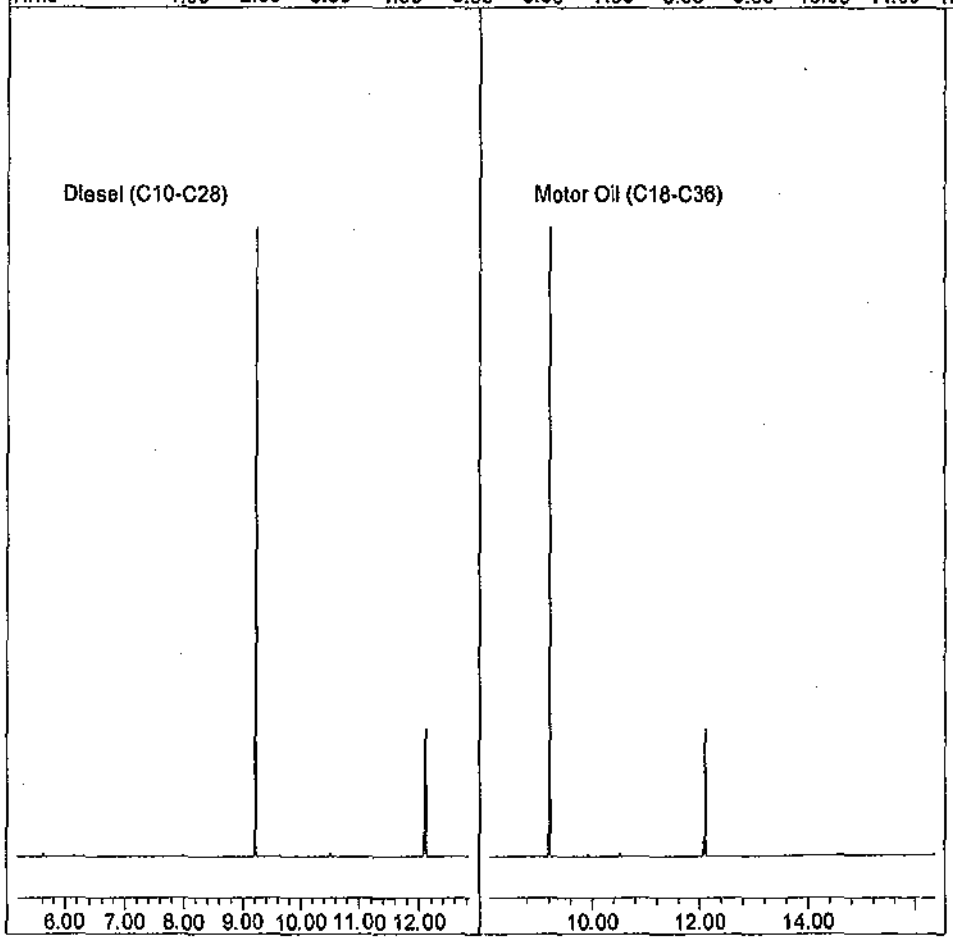
Method : G:\APOLLO\DATA\111115\TPH8S15.M (Chemstation Integrator)
Title : Diesel
Last Update : Thu Nov 17 09:41:49 2011
Response via : Multiple Level Calibration

Volume Inj. : 2UL
Signal Phase : DB-5
Signal Info : FID02A

Compound	R.T.	Response	Conc	Units
System Monitoring Compounds				
4) SC Ortho-Terphenyl(S)	9.21	25946623	40.667	ppb
Surrogate Spike 30.000		Recovery =	135.56%	
6) SC Octacosane(S)	12.11	10118734	43.337	ppb
Surrogate Spike 30.000		Recovery =	144.46%	
Target Compounds				

Data File: G:\APOLLO\DATA\111115\1115025.D
Sample : THC SURR 800/1000





Quantitation Report (Not Reviewed)

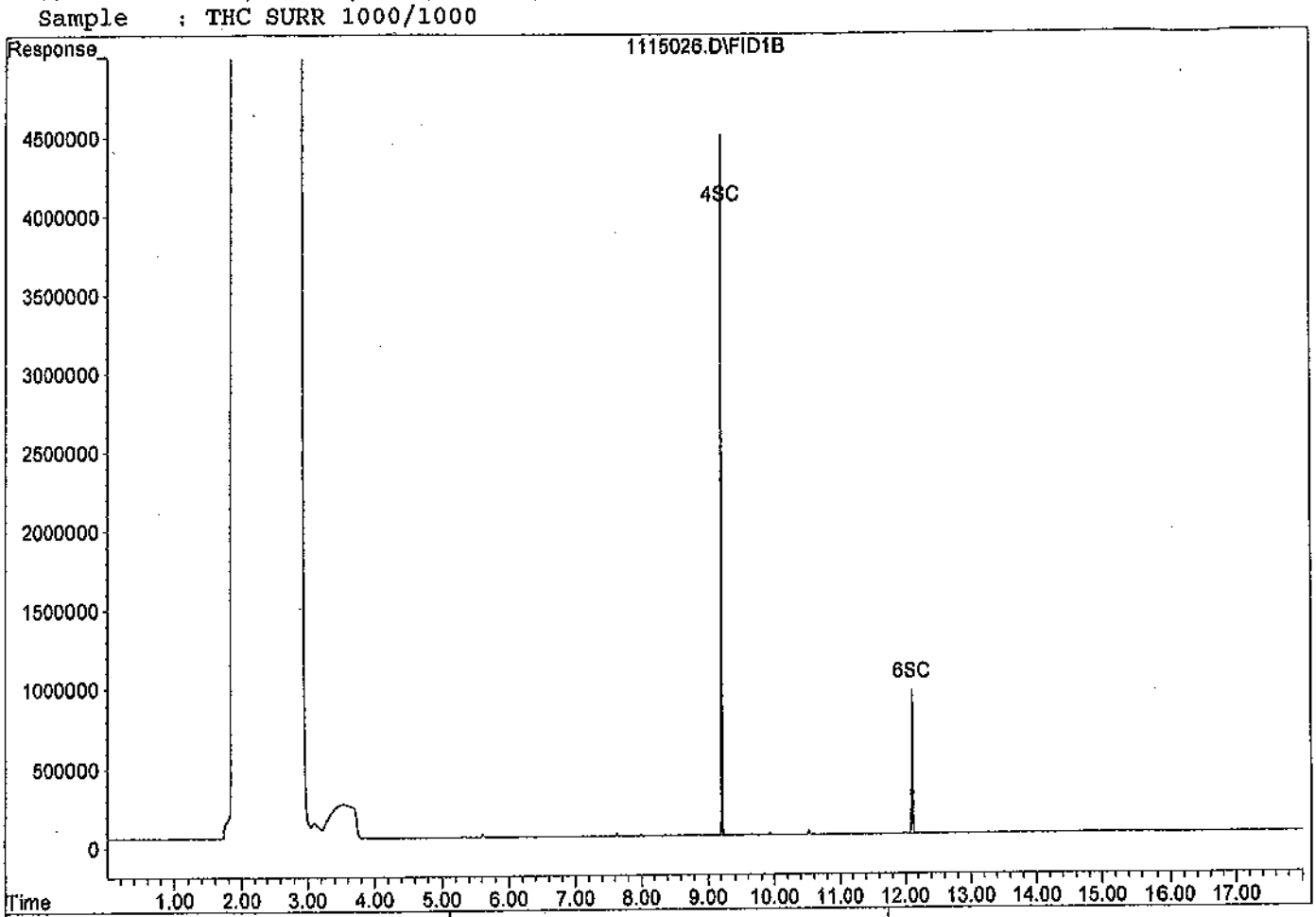
Data File : G:\APOLLO\DATA\111115\1115026.D Vial: 26
Acq On : 11-15-11 20:20:52 Operator: LAC
Sample : THC SURR 1000/1000 Inst : Apollo
Misc : Mix(C) Multiplr: 1.00
IntFile : events.e
Quant Time: Nov 16 14:58 2011 Quant Results File: TPH8S15.RES

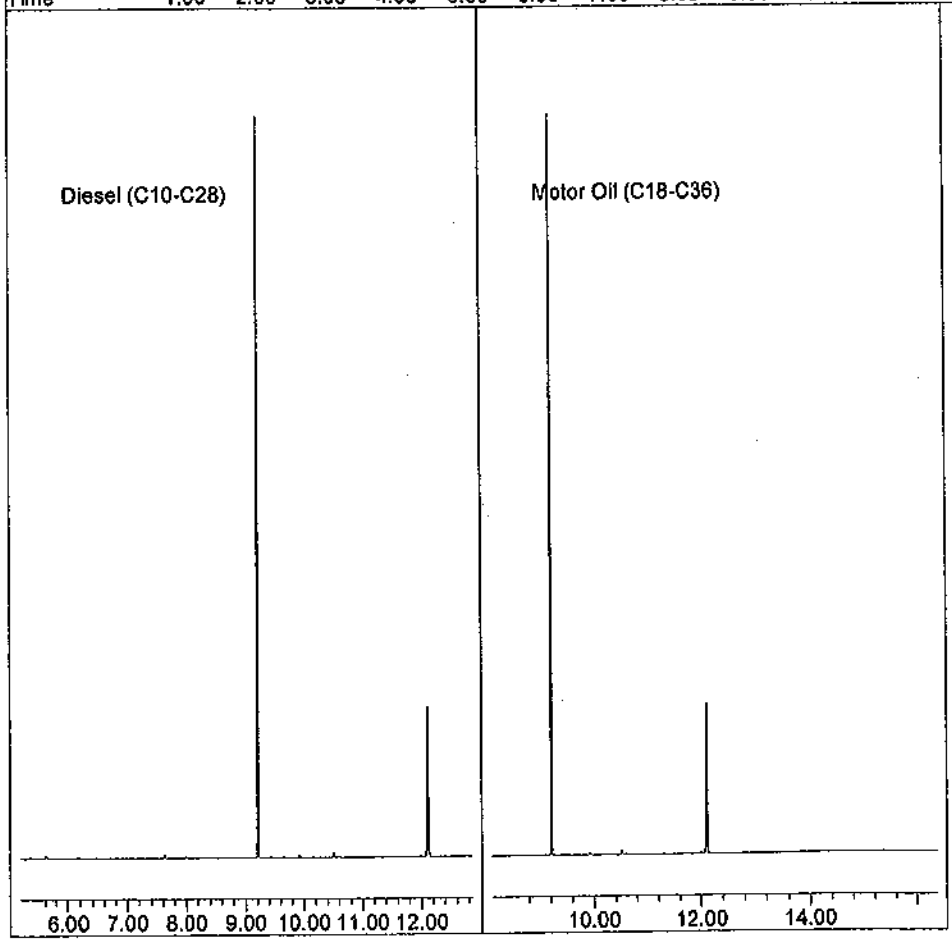
Method : G:\APOLLO\DATA\111115\TPH8S15.M (Chemstation Integrator)
Title : Diesel
Last Update : Thu Nov 17 09:41:49 2011
Response via : Multiple Level Calibration

Volume Inj. : 2UL
Signal Phase : DB-5
Signal Info : FID02A

Compound	R.T.	Response	Conc	Units
System Monitoring Compounds				
4) SC Ortho-Terphenyl(S)	9.21	30736073	48.174	ppb
Surrogate Spike 30.000	Recovery =		160.58%	
6) SC Octacosane(S)	12.11	12029686	51.522	ppb
Surrogate Spike 30.000	Recovery =		171.74%	
Target Compounds				

Data File: G:\APOLLO\DATA\111115\1115026.D
Sample : THC SURR 1000/1000





Quantitation Report (Not Reviewed)

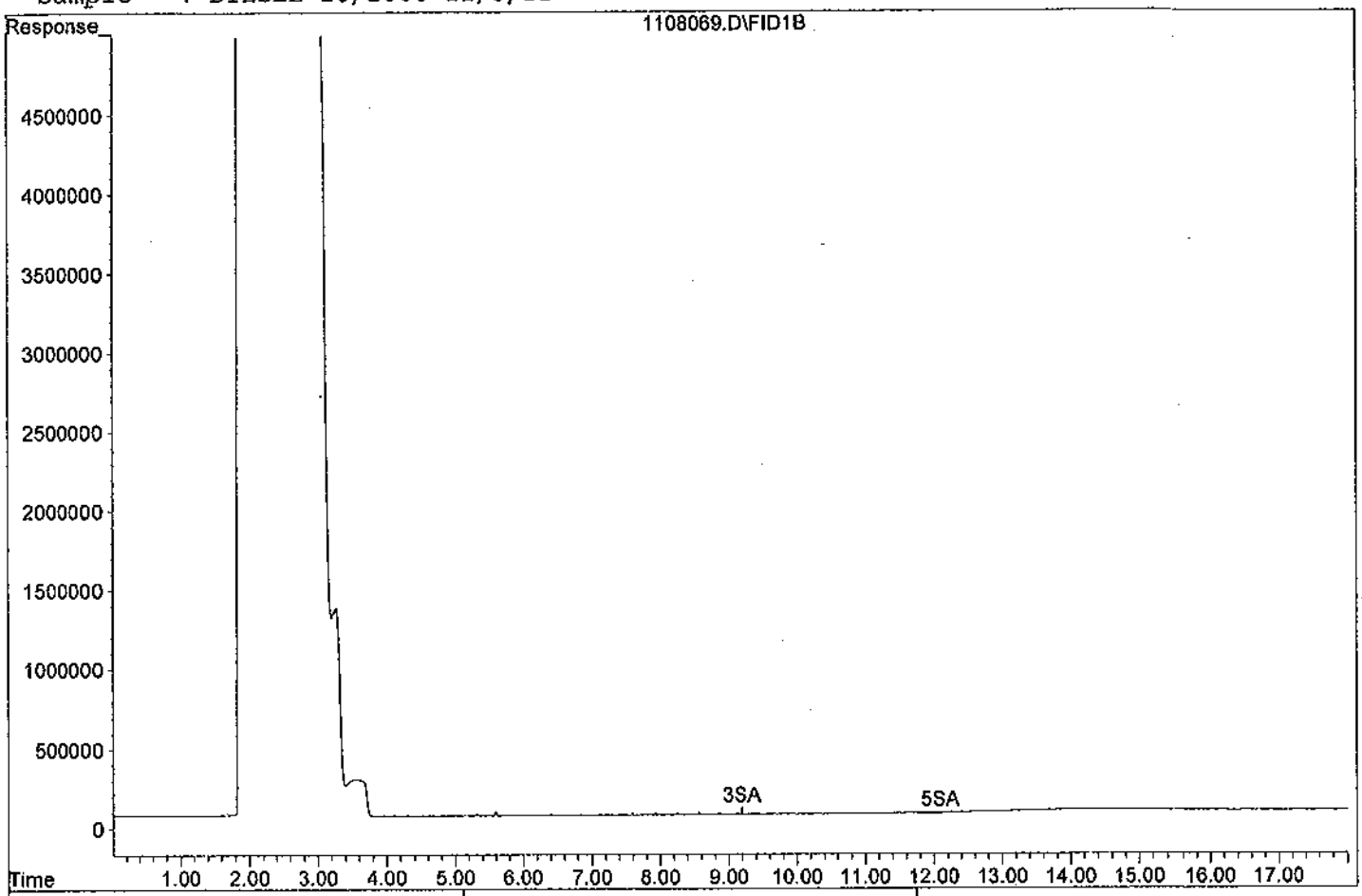
Data File : G:\APOLLO\DATA\111108\1108069.D Vial: 69
Acq On : 11-9-11 17:18:58 Operator: LAC
Sample : DIESEL 10/1000 11/8/11 Inst : Apollo
Misc : Mix(A) Multiplr: 1.00
IntFile : events.e
Quant Time: Nov 16 9:53 2011 Quant Results File: TPH8S15.RES

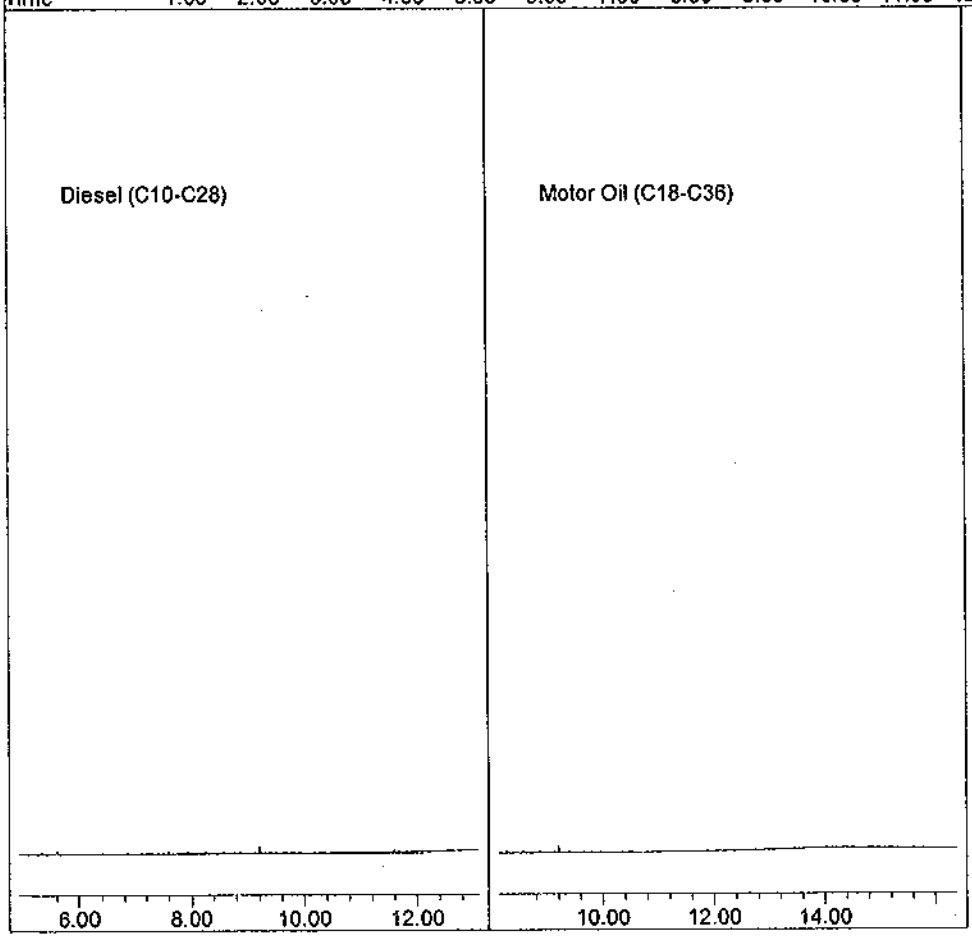
Method : G:\APOLLO\DATA\111108\TPH8S15.M (Chemstation Integrator)
Title : Diesel
Last Update : Wed Nov 16 09:55:03 2011
Response via : Multiple Level Calibration

Volume Inj. : 2UL
Signal Phase : DB-5
Signal Info : FID02A

Compound	R.T.	Response	Conc Units
System Monitoring Compounds			
3) SA Not Used(S)	9.20	302444	0.297 ppb
Surrogate Spike 30.000	Recovery	=	0.99%
5) SA Not Used2(S)	12.10	625179	2.122 ppb
Surrogate Spike 30.000	Recovery	=	7.07%
Target Compounds			
1) HATM Diesel (C10-C28)	9.01	12262633	1055.198 ppb

Data File: G:\APOLLO\DATA\111108\1108069.D
Sample : DIESEL 10/1000 11/8/11





TPH Extractables
TPH8S15

Form 7

Second Source Calibration

Lab Name: APPL, Inc.
Case No:
Matrix:

SDG No: 66102
Date Analyzed: 11/09/11
Instrument: Apollo
Initial Cal. Date: 11/08/11
Data File: 1108070.D

		Compound	MEAN	CCRF	%D		%Drift
1	HATM	Diesel (C10-C28)	305473	234327	23	HATML	4.9
2							
3							
4							
5							
6							
7							
8							
9							
10							
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32							
33							
34							
35							
36							
37							
38							
39							
40							
		Average			23.0		

Quantitation Report (Not Reviewed)

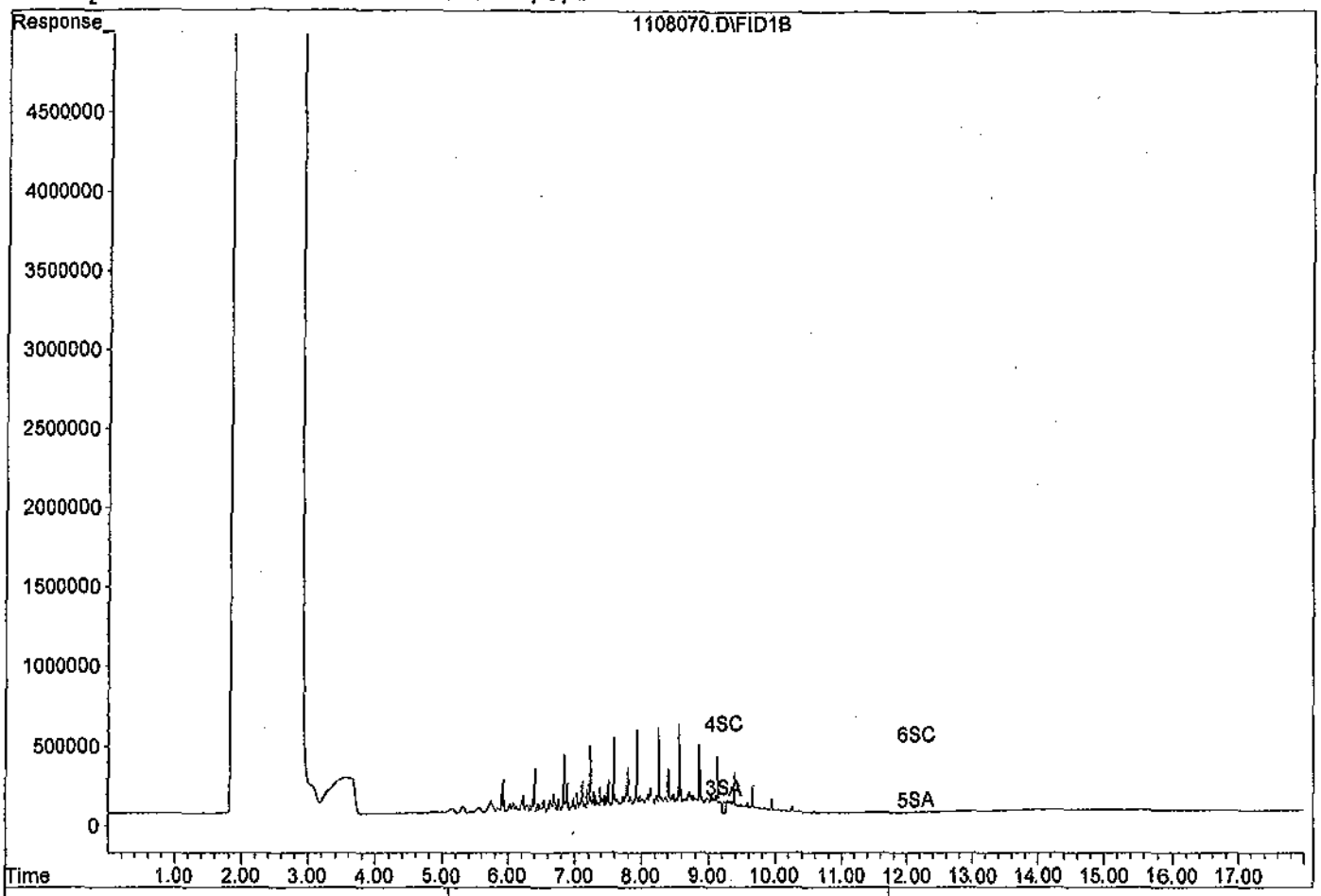
Data File : G:\APOLLO\DATA\111108\1108070.D Vial: 70
Acq On : 11-9-11 17:42:38 Operator: LAC
Sample : DIESEL 400 2ND SRC 11/8/11 Inst : Apollo
Misc : Mix(A) Multiplr: 1.00
IntFile : events.e
Quant Time: Nov 16 9:52 2011 Quant Results File: TPH8S15.RES

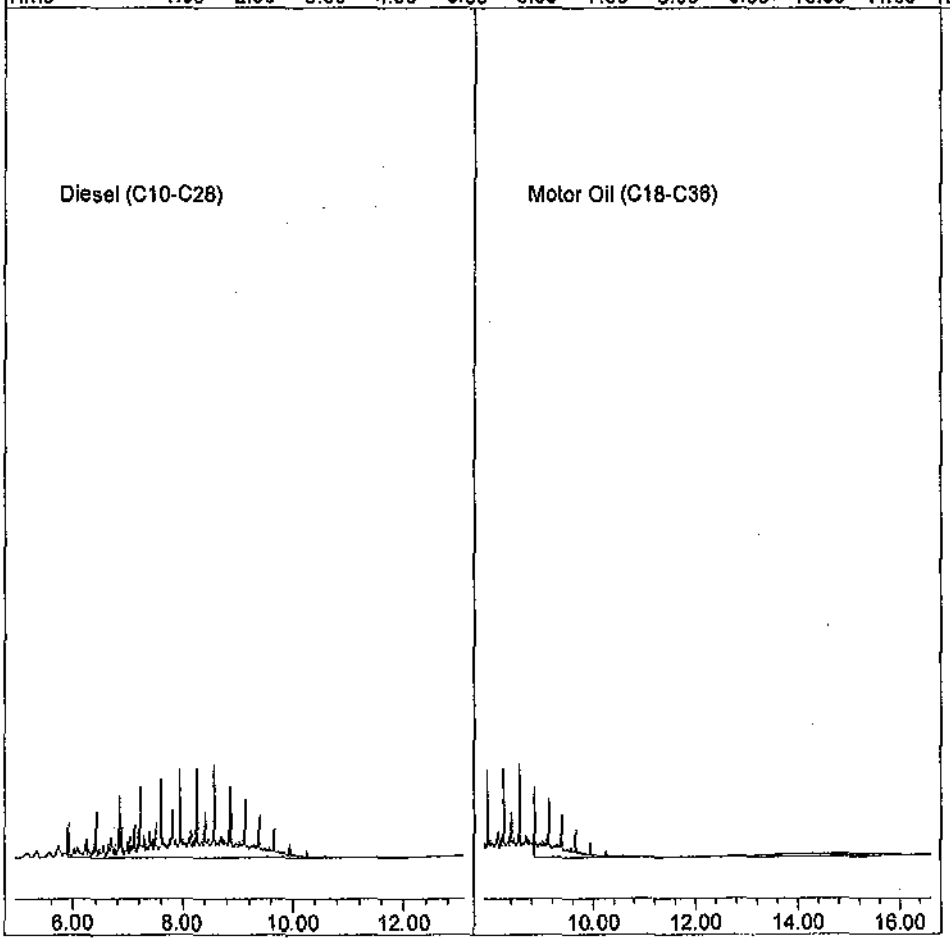
Method : G:\APOLLO\DATA\111115\TPH8S15.M (Chemstation Integrator)
Title : Diesel
Last Update : Thu Nov 17 09:41:49 2011
Response via : Multiple Level Calibration

Volume Inj. : 2UL
Signal Phase : DB-5
Signal Info : FID02A

Compound	R.T.	Response	Conc	Units
System Monitoring Compounds				
3) SA Not Used(S)	9.24	2204277	2.168	ppb
Surrogate Spike 30.000	Recovery =		7.23%	
4) SC Ortho-Terphenyl(S)	9.24	2204277	3.455	ppb
Surrogate Spike 30.000	Recovery =		11.52%	
5) SA Not Used2(S)	12.16	136311	0.463	ppb
Surrogate Spike 30.000	Recovery =		1.54%	
6) SC Octacosane(S)	12.16	136311	0.584	ppb
Surrogate Spike 30.000	Recovery =		1.95%	
Target Compounds				
2) HBTM Motor Oil (C18-C36)	12.24	65481078	788.357	ppb

Data File: G:\APOLLO\DATA\111108\1108070.D
Sample : DIESEL 400 2ND SRC 11/8/11





TPH Extractables
TPH8S15

Form 7

Continuing Calibration

Lab Name: APPL, Inc.
Case No:
Matrix:

SDG No: 66102
Date Analyzed: 11/29/11
Instrument: Apollo
Initial Cal. Date: 11/08/11
Data File: 1129012.D

		Compound	MEAN	CCRF	%D		%Drift
1	HATM	Diesel (C10-C28)	305473	265441	13	HATML	7.9
2							
3							
4							
5							
6							
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36							
37							
38							
39							
40							
		Average			13.0		

Quantitation Report (Not Reviewed)

Data File : G:\APOLLO\DATA\111129\1129012.D Vial: 12
Acq On : 11-29-11 13:06:20 Operator: LAC
Sample : DIESEL 400/1000 11/29/11 Inst : Apollo
Misc : Mix(A) Multiplr: 1.00
IntFile : events.e
Quant Time: Nov 29 16:12 2011 Quant Results File: TPH8S15.RES

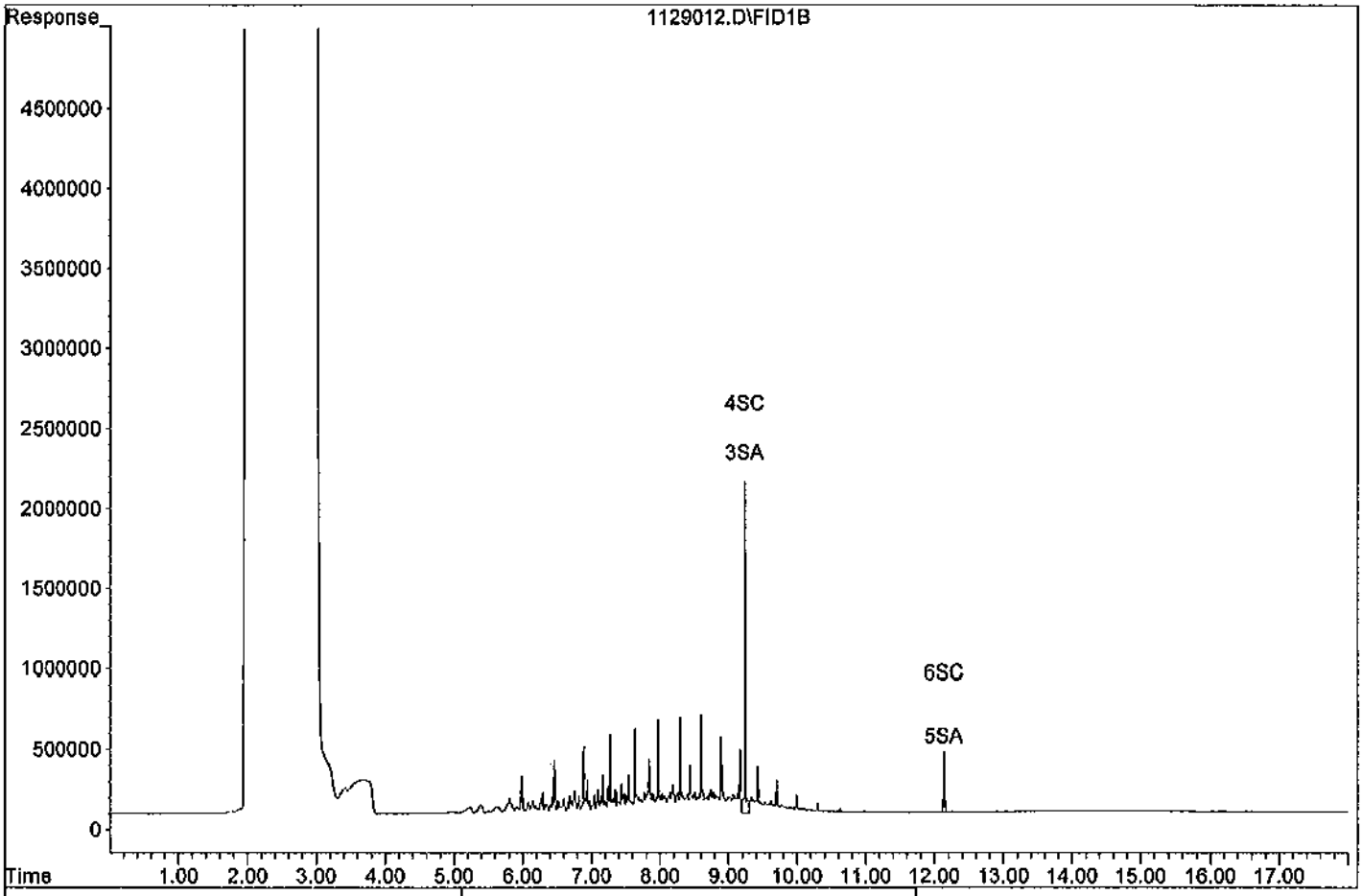
Method : G:\APOLLO\DATA\111123\TPH8S15.M (Chemstation Integrator)
Title : Diesel
Last Update : Mon Nov 28 16:45:51 2011
Response via : Multiple Level Calibration

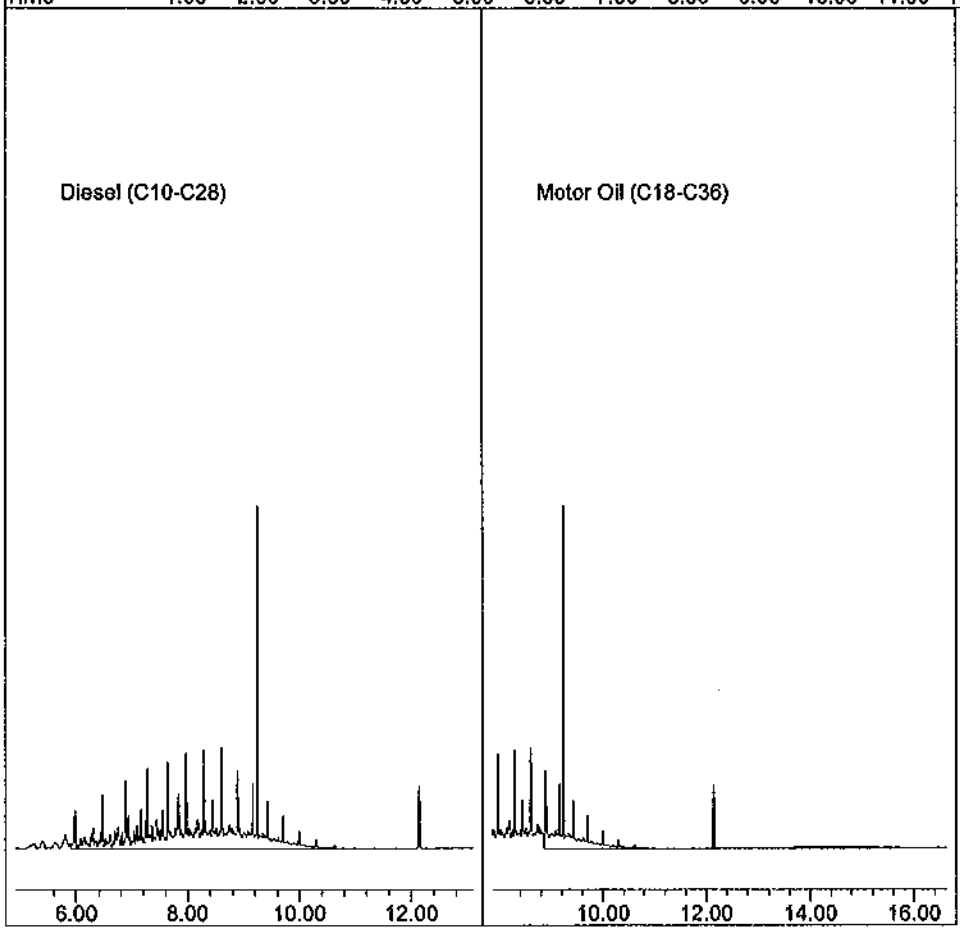
Volume Inj. : 2UL
Signal Phase : DB-5
Signal Info : FID02A

Compound	R.T.	Response	Conc	Units
System Monitoring Compounds				
3) SA Not Used(S)	9.24	18448540	27.180	ppb
Surrogate Spike 30.000	Recovery =		90.60%	
4) SC Ortho-Terphenyl(S)	9.24	18448540	28.915	ppb
Surrogate Spike 30.000	Recovery =		96.38%	
5) SA Not Used2(S)	12.14	5020781	31.843	ppb
Surrogate Spike 30.000	Recovery =		106.14%	
6) SC Octacosane(S)	12.14	5020781	21.038	ppb
Surrogate Spike 30.000	Recovery =		70.13%	
Target Compounds				
1) HATM Diesel (C10-C28)	9.01	212352743	431.747	ppb
2) HBTM Motor Oil (C18-C36)	12.24	65096224	280.338	ppb

(f)=RT Delta > 1/2 Window 127 (m)=manual int.

Data File: G:\APOLLO\DATA\111129\1129012.D
Sample : DIESEL 400/1000 11/29/11





TPH Extractables
TPH8S15

Form 7

Continuing Calibration

Lab Name: APPL, Inc.
Case No:
Matrix:

SDG No: 66102
Date Analyzed: 11/29/11
Instrument: Apollo
Initial Cal. Date: 11/08/11
Data File: 1129024.D

		Compound	MEAN	CCRF	%D	%Drift	
1	HATM	Diesel (C10-C28)	305473	278570	8.8	HATML	13
2							
3							
4							
5							
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34							
35							
36							
37							
38							
39							
40							
		Average			8.8		

Quantitation Report (Not Reviewed)

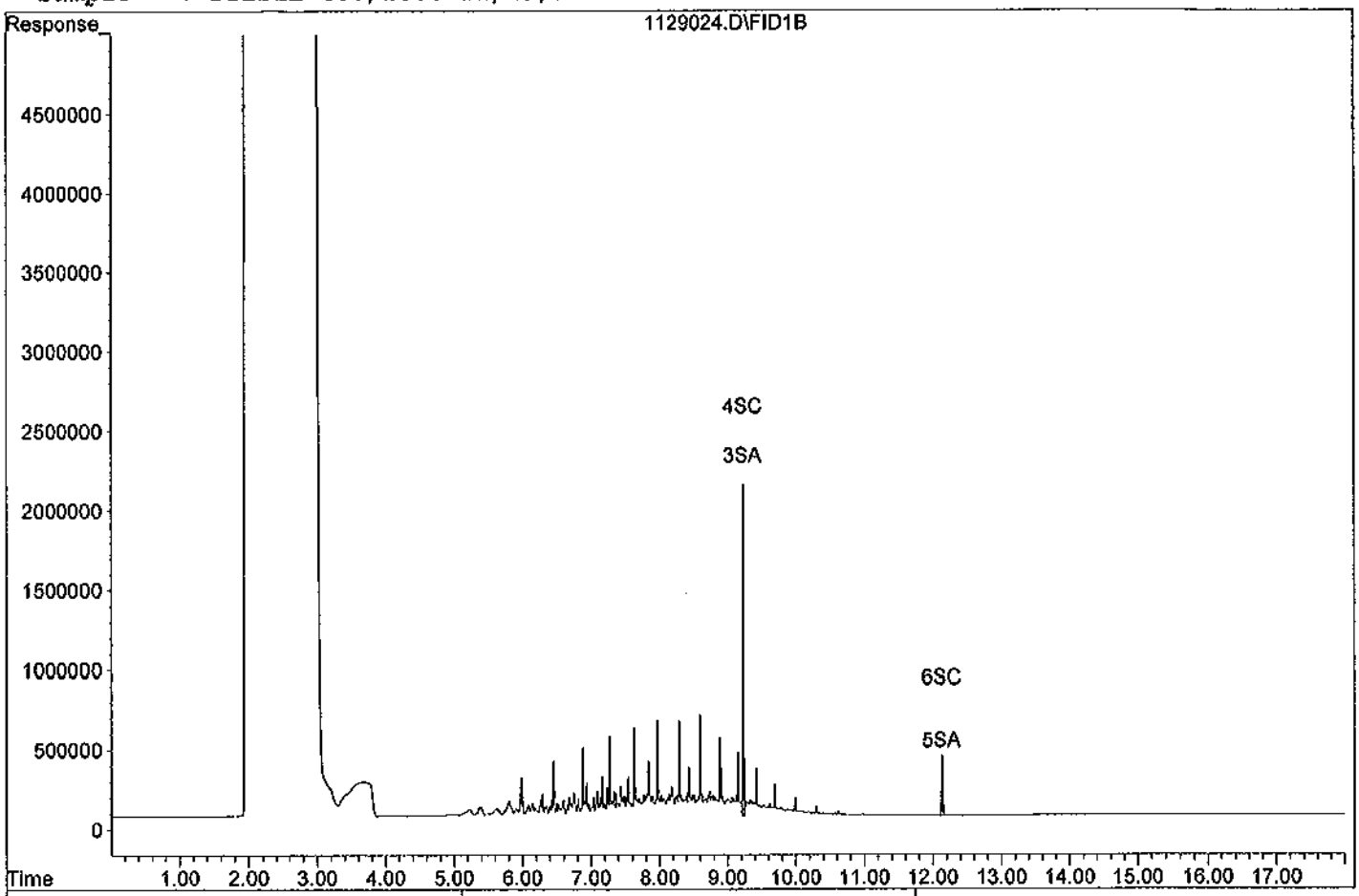
Data File : G:\APOLLO\DATA\111129\1129024.D Vial: 24
Acq On : 11-29-11 21:29:15 Operator: LAC
Sample : DIESEL 400/1000 11/29/11 Inst : Apollo
Misc : Mix(A) Multiplr: 1.00
IntFile : events.e
Quant Time: Nov 30 8:35 2011 Quant Results File: TPH8S15.RES

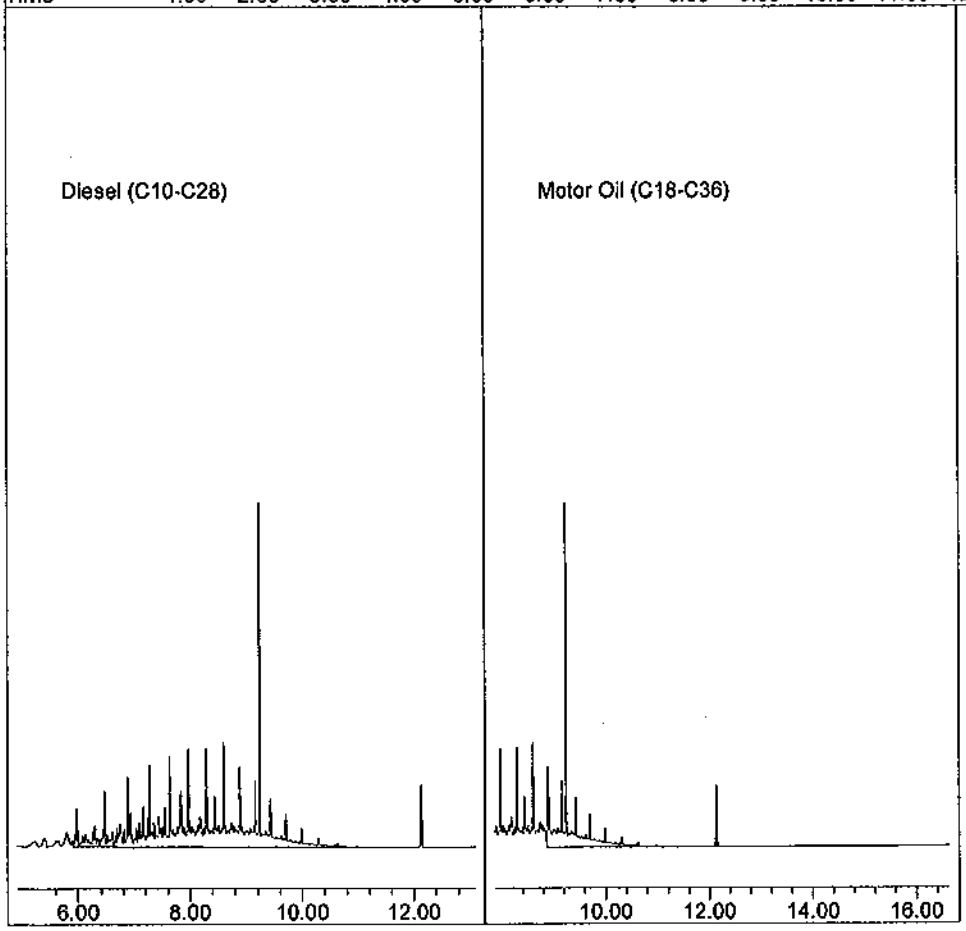
Method : G:\APOLLO\DATA\111123\TPH8S15.M (Chemstation Integrator)
Title : Diesel
Last Update : Mon Nov 28 16:45:51 2011
Response via : Multiple Level Calibration

Volume Inj. : 2UL
Signal Phase : DB-5
Signal Info : FID02A

Compound	R.T.	Response	Conc	Units
System Monitoring Compounds				
3) SA Not Used(S)	9.24	15158305	22.332	ppb
Surrogate Spike 30.000	Recovery =		74.44%	
4) SC Ortho-Terphenyl(S)	9.24	15158305	23.758	ppb
Surrogate Spike 30.000	Recovery =		79.19%	
5) SA Not Used2(S)	12.13	4847180	30.742	ppb
Surrogate Spike 30.000	Recovery =		102.47%	
6) SC Octacosane(S)	12.13	4847180	20.311	ppb
Surrogate Spike 30.000	Recovery =		67.70%	
Target Compounds				
1) HATM Diesel (C10-C28)	9.01	222856237	453.399	ppb
2) HBTM Motor Oil (C18-C36)	12.24	69450604	299.090	ppb

Data File: G:\APOLLO\DATA\111129\1129024.D
Sample : DIESEL 400/1000 11/29/11





EPA 8015 Modified
Total Petroleum Hydrocarbons
Raw Data

Method Blank
TPH Diesel Water

APPL Inc.
908 North Temperance Avenu
Clovis, CA 93611

Blank Name/QCG: **111031W-49334 - 160886**
Batch ID: #TPETD-111031A

Sample Type	Analyte	Result	LOQ	LOD	DL	Units	Extraction Date	Analysis Date
BLANK	DIESEL FUEL	80.8 U	150	80.8	40.4	ug/L	10/31/11	11/06/11
BLANK	SURROGATE: OCTACOSANE (S)	71.2	28-142			%	10/31/11	11/06/11
BLANK	SURROGATE: ORTHO-TERPHEN	60.5	57-132			%	10/31/11	11/06/11

Quant Method: TPH1028.M
Run #: 1106005
Instrument: Apollo
Sequence: 111106
Initials: LA

GC SC-Blank-REG MDLs
Printed: 11/30/11 11:27:27 AM

Quantitation Report (QT Reviewed)

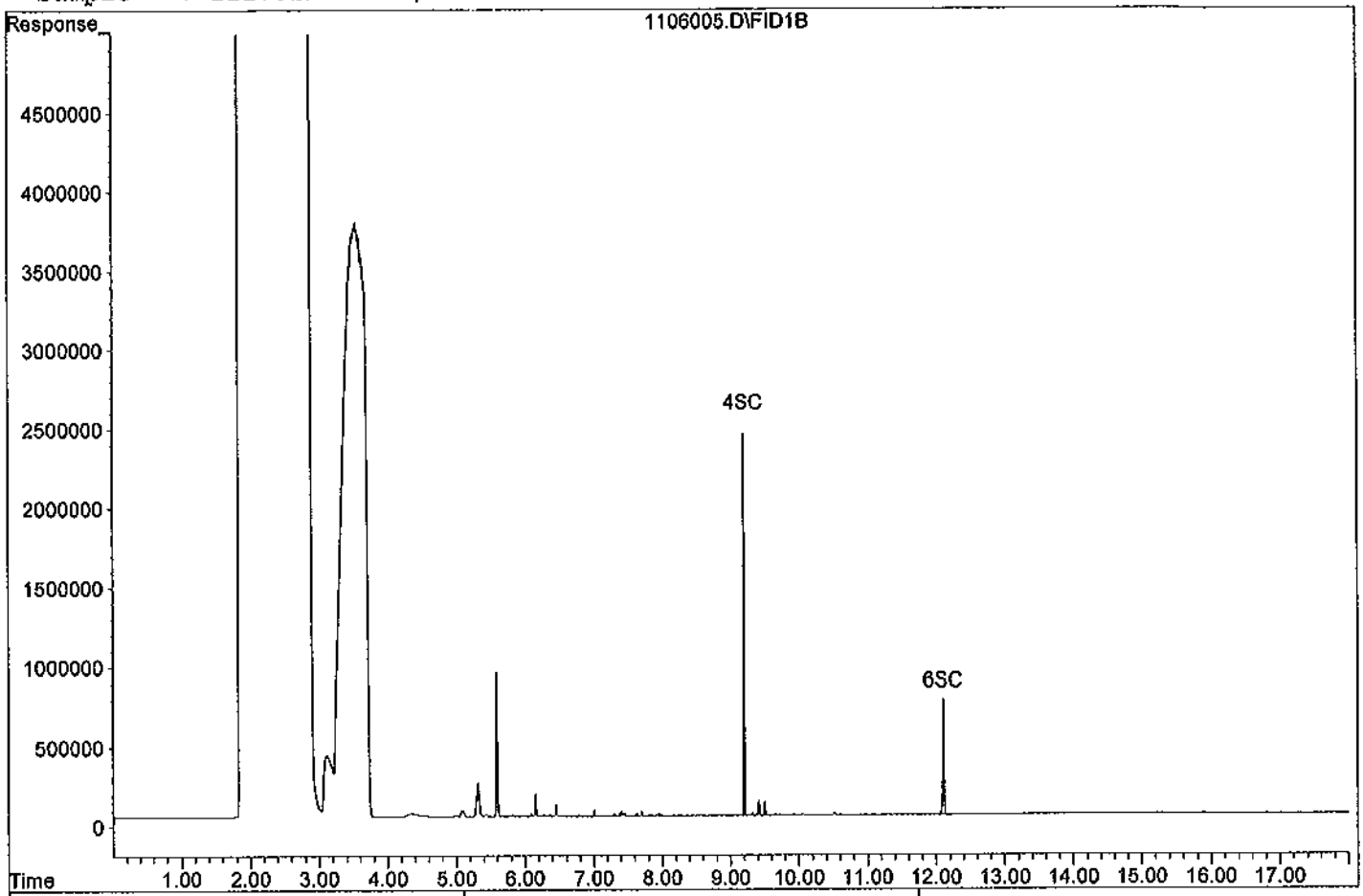
Data File : G:\APOLLO\DATA\111106\1106005.D Vial: 5
Acq On : 11-6-11 17:22:09 Operator: LAC
Sample : 111031A BLK 5/1000 Inst : Apollo
Misc : Water Multiplr: 5.00
IntFile : events.e
Quant Time: Nov 7 9:44 2011 Quant Results File: TPH1028.RES

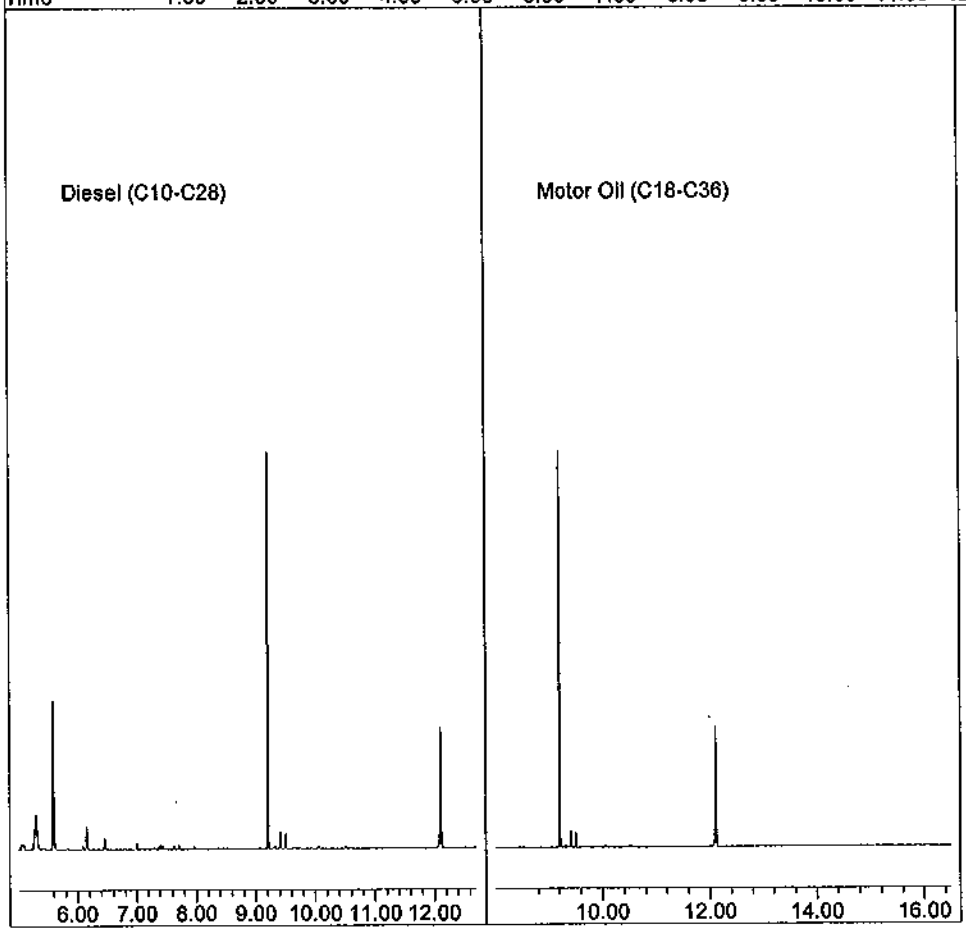
Method : G:\APOLLO\DATA\111028\TPH1028.M (Chemstation Integrator)
Title : Diesel
Last Update : Mon Oct 31 10:02:11 2011
Response via : Multiple Level Calibration

Volume Inj. : 2UL
Signal Phase : DB-5
Signal Info : FID02A

Compound	R.T.	Response	Conc Units
System Monitoring Compounds			
4) SC Ortho-Terphenyl(S)	9.21	15818444	90.703 ppb
Surrogate Spike 150.000		Recovery =	60.47%
6) SC Octacosane(S)	12.11	9943034	106.834 ppb
Surrogate Spike 150.000		Recovery =	71.22%
Target Compounds			

Data File: G:\APOLLO\DATA\111106\1106005.D
Sample : 111031A BLK 5/1000





Laboratory Control Spike Recovery
TPH Diesel Water

APPL ID: **111031W-49334 LCS - 160886**
Batch ID: #TPETD-111031A

APPL Inc.
908 North Temperance Avenue
Clovis, CA 93611

Compound Name	Spike Level ug/L	SPK Result ug/L	SPK % Recovery	Recovery Limits
DIESEL FUEL	2000	1520	76.0	61-143
SURROGATE: OCTACOSANE (S)	150	125	83.3	28-142
SURROGATE: ORTHO-TERPHENYL (S)	150	147	98.0	57-132

Comments:

Primary	SPK
Quant Method :	TPH8S15.M
Extraction Date :	10/31/11
Analysis Date :	11/29/11
Instrument :	Apollo
Run :	1129017
Initials :	LA

Printed: 11/30/11 11:27:21 AM

APPL Standard LCS

Quantitation Report (QT Reviewed)

Data File : G:\APOLLO\DATA\111129\1129017.D Vial: 17
Acq On : 11-29-11 18:45:15 Operator: LAC
Sample : 111031A LCS-1 5/1000 Inst : Apollo
Misc : Water Multiplr: 5.00
IntFile : events.e
Quant Time: Nov 30 10:29 2011 Quant Results File: TPH8S15.RES

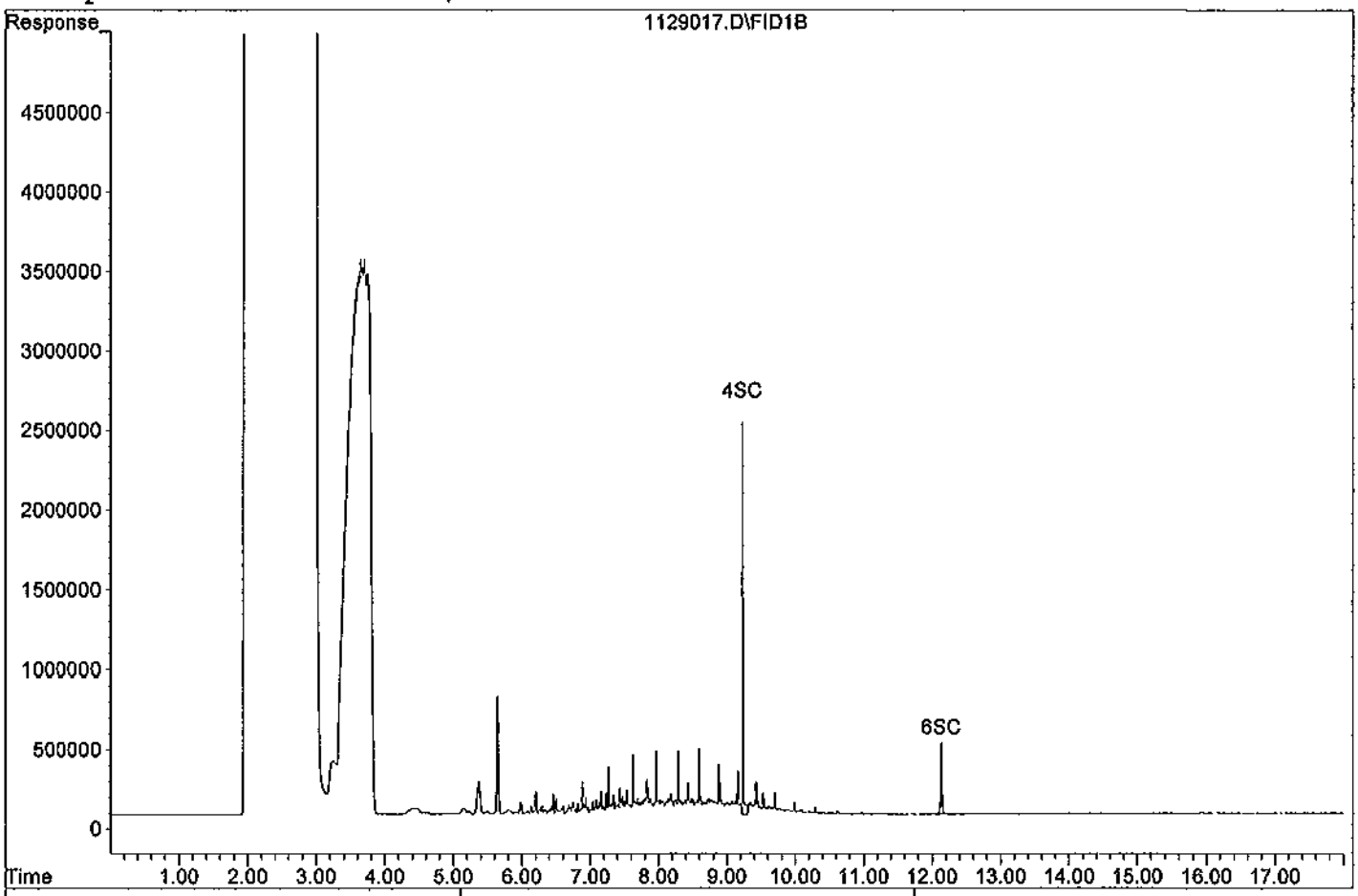
Method : G:\APOLLO\DATA\111123\TPH8S15.M (Chemstation Integrator)
Title : Diesel
Last Update : Mon Nov 28 16:45:51 2011
Response via : Multiple Level Calibration

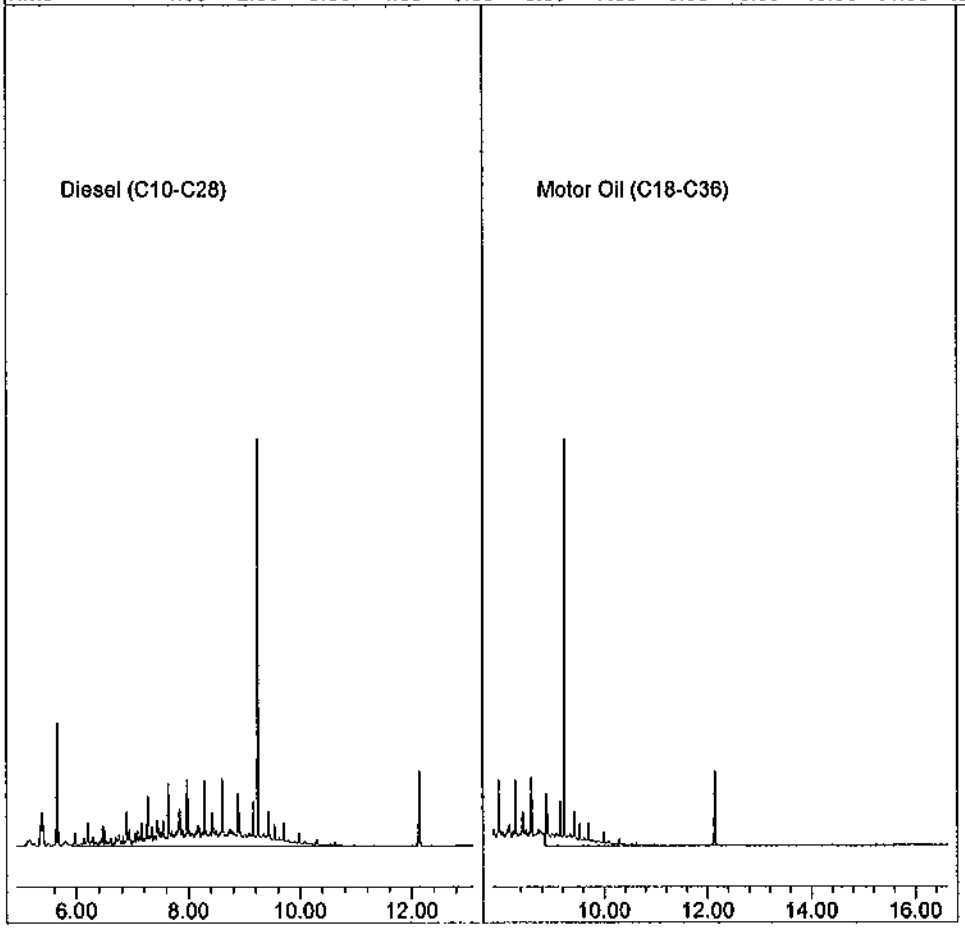
Volume Inj. : 2UL
Signal Phase : DB-5
Signal Info : FID02A

Compound	R.T.	Response	Conc Units
System Monitoring Compounds			
4) SC Ortho-Terphenyl(S)	9.24	18706073	146.595 ppb
Surrogate Spike 150.000		Recovery =	97.73%
6) SC Octacosane(S)	12.14	5944718	124.549 ppb
Surrogate Spike 150.000		Recovery =	83.03%
Target Compounds			
1) HATM Diesel (C10-C28)	9.01	150395375	1520.130 ppb
2) HBTM Motor Oil (C18-C36)	12.24	50801855	1093.895 ppb

Algorithm Check: $\frac{(18706073)(5)}{(319010)(2)} = 146.594\,57227$ LAC 11/30/11 LAC 11/30/11

Data File: G:\APOLLO\DATA\111129\1129017.D
Sample : 111031A LCS-1 5/1000





Matrix Spike Recoveries
TPH Diesel Water

APPL ID: **111031W-49334 MS - 160886**
Batch ID: #TPETD-111031A
Sample ID: AY49334
Client ID: ES047

APPL Inc.
908 North Temperance Avenue
Clovis, CA 93611

Compound Name	Spike Lvl ug/L	Matrix Result ug/L	SPK Result ug/L	DUP Result ug/L	SPK % Recovery	DUP % Recovery	Recovery Limits	RPD %	RPD Limits
DIESEL FUEL	2000	730	1400	1330	33.5 #	30.0 #	61-143	5.1	30
SURROGATE: OCTACOSANE (S)	150	NA	131	125	87.3	83.3	28-142		
SURROGATE: ORTHO-TERPHENYL (S)	150	NA	108	89.3	72.0	59.5	57-132		

= Recovery is outside QC limits.

Comments:

Primary	SPK	DUP
Quant Method :	TPH1028.M	TPH1028.M
Extraction Date :	10/31/11	10/31/11
Analysis Date :	11/06/11	11/06/11
Instrument :	Apollo	Apollo
Run :	1106014	1106015
Initials :	LA	

Printed: 11/30/11 11:27:18 AM
APPL MSD SCII

Quantitation Report (QT Reviewed)

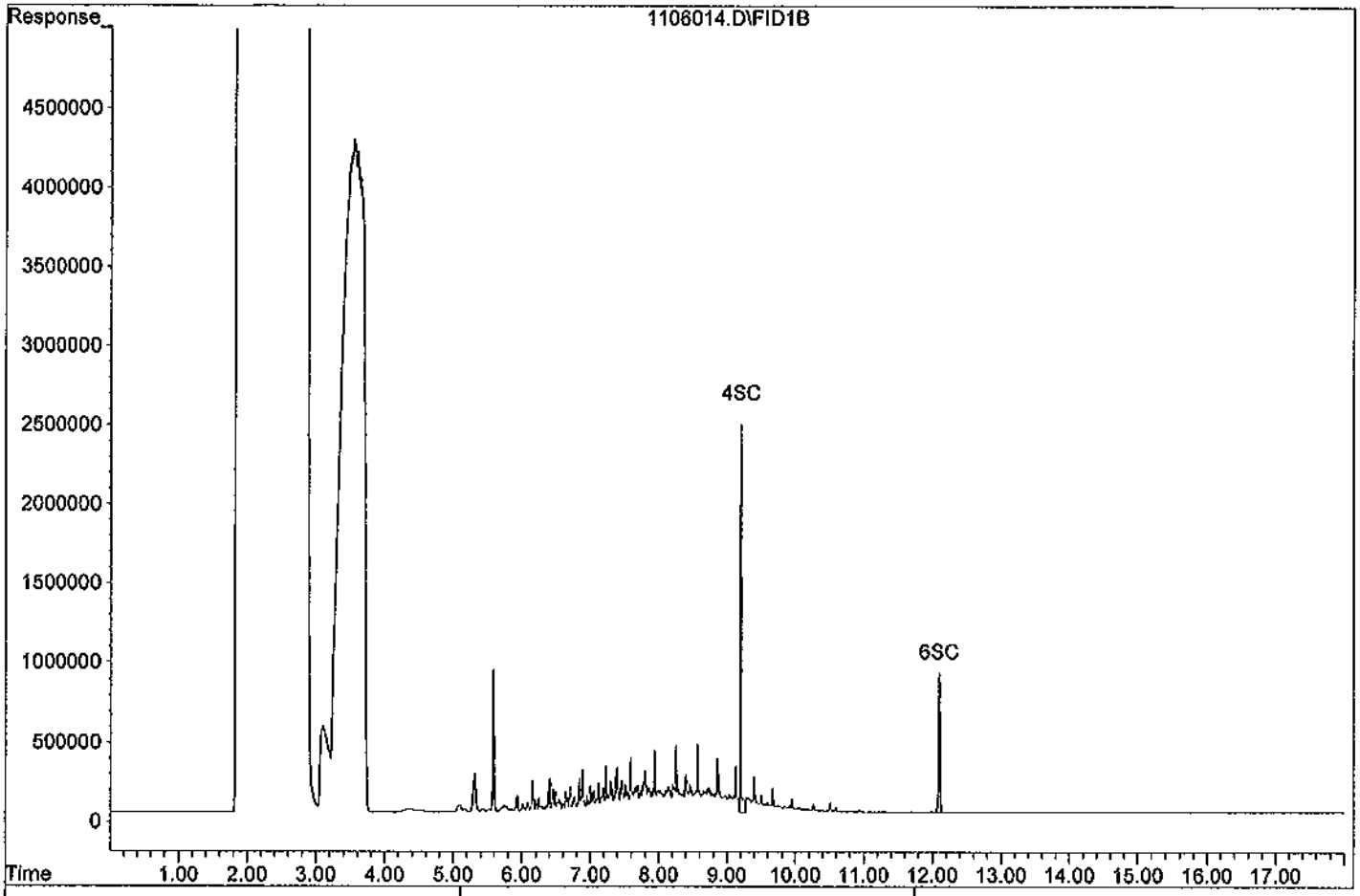
Data File : G:\APOLLO\DATA\111106\1106014.D Vial: 14
Acq On : 11-6-11 20:53:41 Operator: LAC
Sample : AY49334W35 MS-1 5/1030 Inst : Apollo
Misc : Water Multiplr: 4.85
IntFile : events.e
Quant Time: Nov 7 9:50 2011 Quant Results File: TPH1028.RES

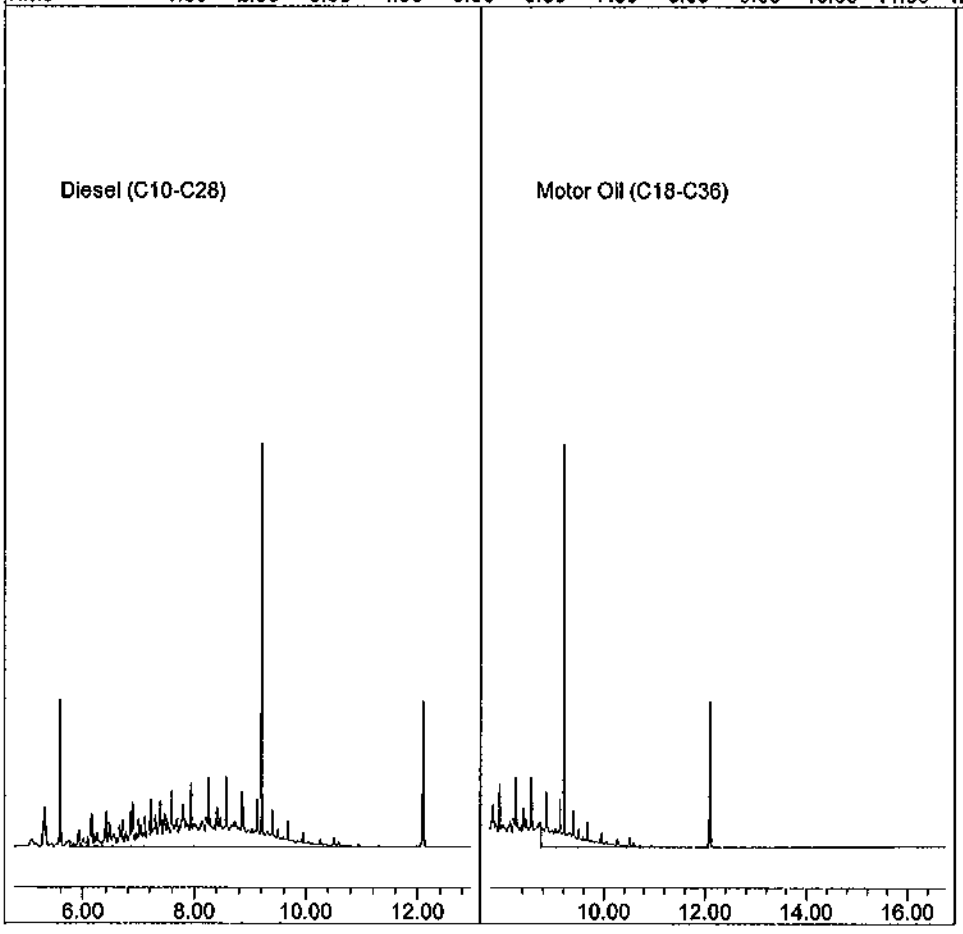
Method : G:\APOLLO\DATA\111028\TPH1028.M (Chemstation Integrator)
Title : Diesel
Last Update : Mon Oct 31 10:02:11 2011
Response via : Multiple Level Calibration

Volume Inj. : 2UL
Signal Phase : DB-5
Signal Info : FID02A

Compound	R.T.	Response	Conc Units
System Monitoring Compounds			
4) SC Ortho-Terphenyl(S)	9.21	19333669	107.631 ppb
Surrogate Spike 145.631		Recovery =	73.91%
6) SC Octacosane(S)	12.10	12570361	131.129 ppb
Surrogate Spike 145.631		Recovery =	90.04%
Target Compounds			
1) HATM Diesel (C10-C28)	8.86	243258357	1402.635 ppb
2) HBTM Motor Oil (C18-C36)	12.25	67181515	899.542 ppb

Data File: G:\APOLLO\DATA\111106\1106014.D
Sample : AY49334W35 MS-1 5/1030





Quantitation Report (QT Reviewed)

Data File : G:\APOLLO\DATA\111106\1106015.D Vial: 15
Acq On : 11-6-11 21:17:05 Operator: LAC
Sample : AY49334W36 MSD-1 5/1030 Inst : Apollo
Misc : Water Multiplr: 4.85
IntFile : events.e
Quant Time: Nov 7 9:50 2011 Quant Results File: TPH1028.RES

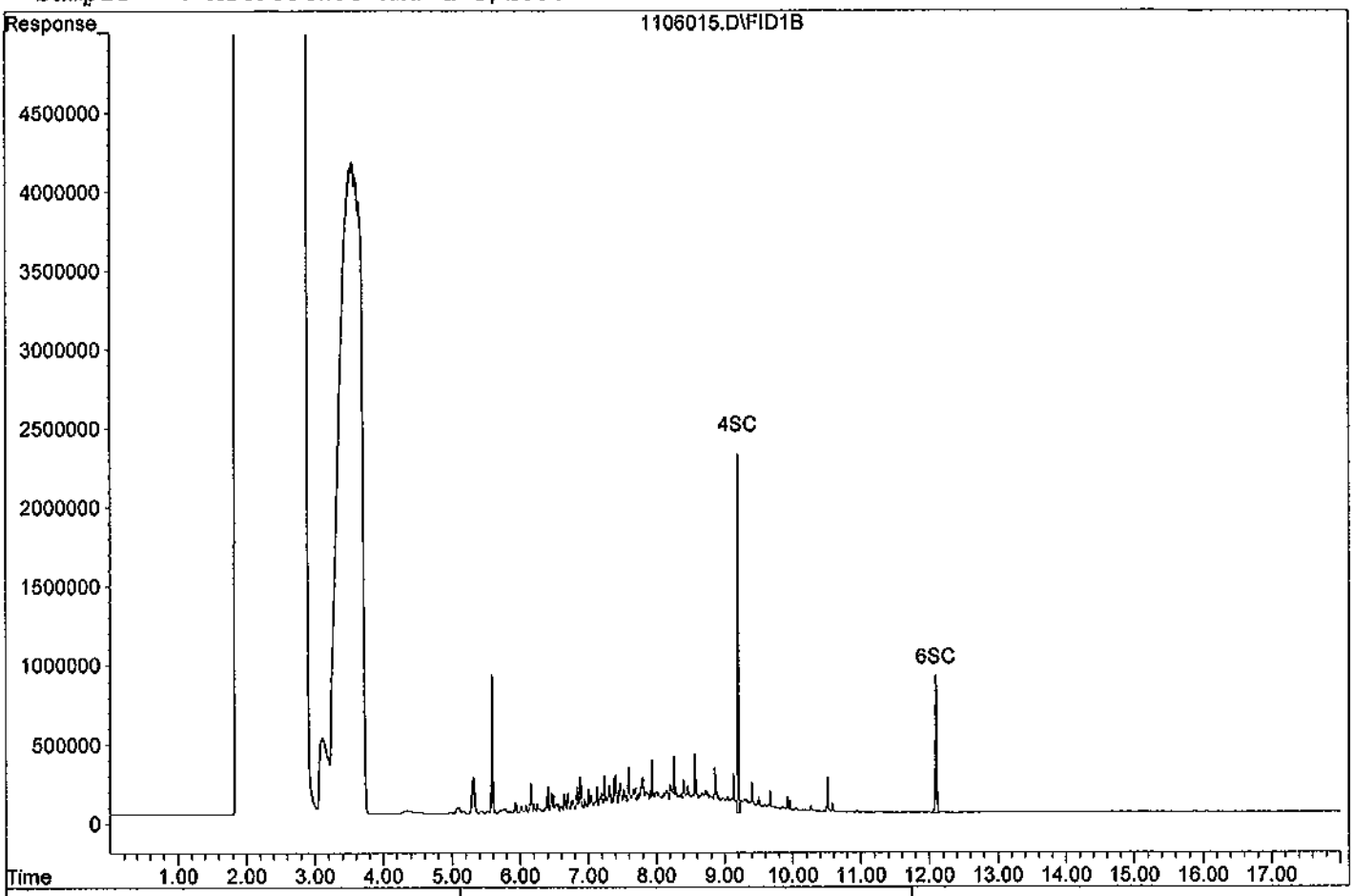
Method : G:\APOLLO\DATA\111028\TPH1028.M (Chemstation Integrator)
Title : Diesel
Last Update : Mon Oct 31 10:02:11 2011
Response via : Multiple Level Calibration

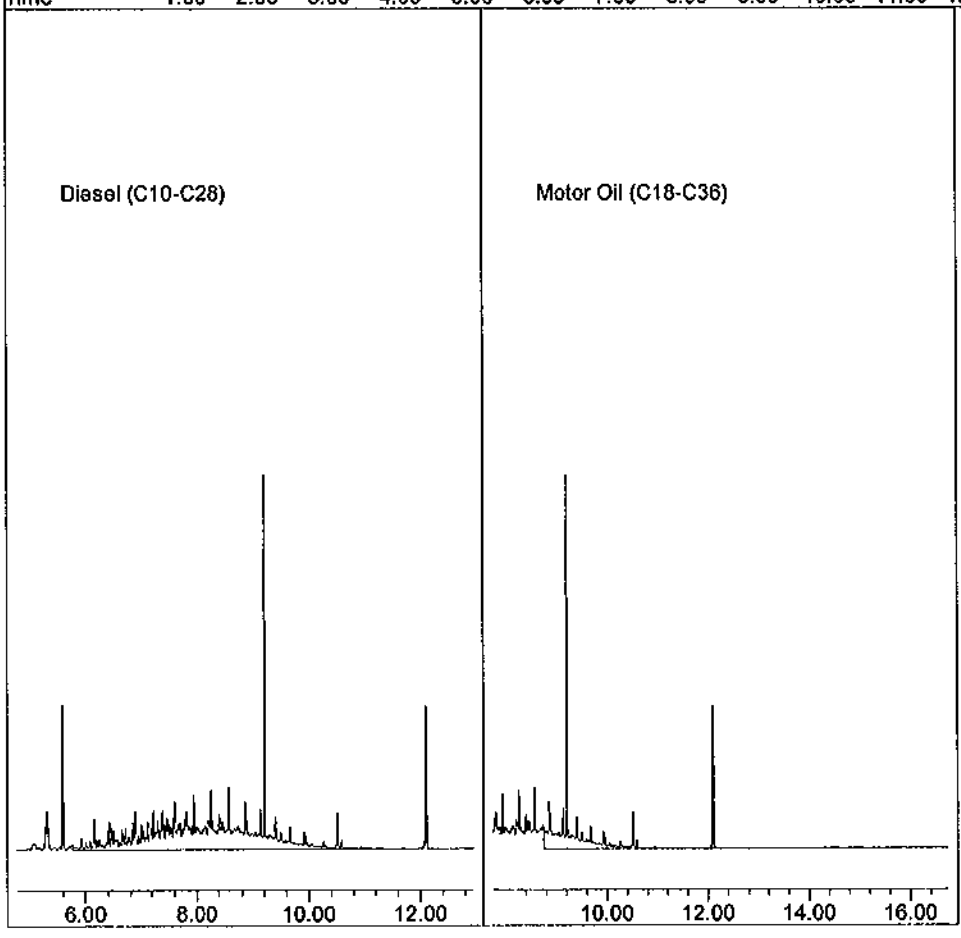
Volume Inj. : 2UL
Signal Phase : DB-5
Signal Info : FID02A

Compound	R.T.	Response	Conc Units
System Monitoring Compounds			
4) SC Ortho-Terphenyl(S)	9.21	16036435	89.275 ppb
Surrogate Spike 145.631		Recovery =	61.30%
6) SC Octacosane(S)	12.10	11966186	124.827 ppb
Surrogate Spike 145.631		Recovery =	85.71%
Target Compounds			
1) HATM Diesel (C10-C28)	8.86	230106893	1326.804 ppb
2) HBTM Motor Oil (C18-C36)	12.25	67186331	899.607 ppb

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

Data File: G:\APOLLO\DATA\111106\1106015.D
Sample : AY49334W36 MSD-1 5/1030





STANDARD	INITIAL CONC	SOURCE	DATE	ALIQUOT	FINAL VOLUME	FINAL CONC	SOLVENT LOT #

029

MOTOR OIL SPIKE

MOTOR OIL — 50,000ug/ml | 0251 | 200ml | 50ML | 2000ug/ml | MC | #070111B | 8/5/11 | EX: 11/5/11

Motor Oil Composite, 50,000 mg/L, 1 ml
o2si smart solutions 116390-02 Storage: </= -10 Degrees C
Made in USA Lot No: 161898 Solvent: Methylene Chloride
Exp: 7/23/2013
Date Op: Motor Oil Composite
Lot #: 161898 - 28613
Rec: 4/14/11 MFR exp. 07/23/13

Motor Oil Composite, 50,000 mg/L, 1 ml
o2si smart solutions 116390-02 Storage: </= -10 Degrees C
Made in USA Lot No: 161898 Solvent: Methylene Chloride
Exp: 7/23/2013
Dat Motor Oil Composite
Lot #: 161898 - 28614
Rec: 4/14/11 MFR exp. 07/23/13

OCL/OP WATER SURROGATE

DBC, TCMX, DECA — 500ug/ml | 0251 | 30ml | 100ML | 1.5ug/ml | ACETONE | #011011C | 8/5/11 | EX: 11/5/11

Pesticide Surrogate Solution, 5,000 mg/L, 1 ml
o2si smart solutions
Cat. No: 130070-02 Exp: 12/19/2012
Lot No: 154164 -27593 Storage: </= Ambient
Solvent: Tol.:Hex. 1:1
Not for Human Consumption For Research Use Only
Made in USA Date Opened: 8/5/11 EX: 8/5/12

TBP, TPP — 100ug/ml | 0251 | 500ml | 5ug/ml

Tributyl- and Triphenylphosphate Soltuion, 1,000 mg/L, 1 ml
o2si smart solutions
Phone: (843) 763-4884 Fax: (843) 766-9182
Cat. No: 130161-02 Exp: 7/12/2012
Lot No: 148444 -27667 Storage: <10 Degrees C
Solvent: Acetone
Not for Human Consumption For Research Use Only
Made in USA Date Opened: 8/5/11 EX: 8/5/12

OCL/OP SOIL SURROGATE

DBC, DECA, TCMX — 500ug/ml | 0251 | 400ml | 100ML | 20ug/ml | ACETONE | #011011C | 8/5/11 | EX: 11/5/11
CAT: 130070-02
LOT: 154164-27593
OP: 8/5/11 EX: 8/5/12

TBP, TPP — 100ug/ml | 0251 | 500ml | 5ug/ml
CAT: 130161-02
LOT: 148444-27667
OP: 8/5/11
EX: 7/12/12

STANDARD	INITIAL CONC	SOURCE DATE	ALIQUOT	FINAL VOLUME	FINAL CONC	SOLVENT LOT #	
		THC SURROGATE			(* GAVE TO EXTRACTION)		
O TERPHENYL	600ug/ml	0251	N/A	25mL	600ug/ml	N/A	8/8/11
OCTACOSANE		CAT: 110316-05					EX: 8/8/12
		LOT: 170258-29333					
		OP: 8/8/11					
		EX: 8/8/12					
		THC SURROGATE					
O TERPHENYL	600ug/ml	0251	N/A	25mL	600ug/ml	N/A	8/8/11
OCTACOSANE		CAT: 110316-05					EX:
	8/8/11	LOT: ~~1704~~ 176405-29334					8/8/12
		THRU					
		176405-29337					
		OP: 8/8/11					
		EX: 8/8/12					
		OCl Degradation Check					
DDT	100ug/ml	0251	250ul	50mL	0.5ug/ml	HEXANE	
DDD		CAT: 130109-01				# 082610B	8/11/11
DDE		LOT: OC Pesticide Degradation Check Lot #: 176400-29311					EX: 2/11/12
ENDRIN		Rec: 8/8/11 MFR exp. 07/29/14					
		Op: 8/11/11	8/11/11				
ENDRIN KETONE		EX: 8/11/12					
ENDRIN ALDEHYDE							
		OCL STOCK					
VARIOUS	100ug/ml	0251	1000ul	10mL	10ug/ml	Hexane	
Analytes		Organochlorine Pesticide Solution 20, 100 mg/L, 1ml				#082610B	8/11/11
		130615-09					EX: 8/11/12
		Lot # 176673 Storage ≤6 Degrees C Expiry 8/2/14	8/11/11				
		Organochlorine Pest Soln 20	OP: 8/11/11				
		Lot #: 176673-29347	EX: 8/11/12				
		Rec: 8/9/11 MFR exp. 08/02/14					

STANDARD 052	INITIAL CONC	SOURCE DATE	ALIQUOT	FINAL VOLUME	FINAL CONC	SOLVENT LOT #	DATE / INITIALS
DIESEL STANDARD							
DIESEL FUEL #2	50,000mg/L	0251	1000mL	50mL	1000mg/L	MC # 051711B	9/1/11 EX: 3/1/12
	Diesel Fuel #2 Composite, 50,000 mg/L, 1 ml 911598-83 Lot # 167768 Storage 5-10 Degrees C Expiry 2/15/15 Made in the USA Solvent: Methylene Chloride Diesel Fuel #2 Composite OP: 9/1/11 Lot #: 167768 - 28176 EX: 9/1/12 Rec: 1/20/11 MFR exp. 02/15/15						
OCTACOSANE	1000mg/L	0251	4170mL		50mg/L		
O-TERPHENYL	CAT: 110316-05 LOT: 176405-29337 OP: 9/1/11 EX: 9/1/12						
MOTOR OIL STANDARD							
MOTOR OIL	50,000mg/L	0251	1000mL	50mL	1000mg/L	MC # 051711B	9/1/11 EX: 3/1/12
	O2si SMART SOLUTIONS 116390-02 Motor Oil Composite, 50,000 mg/L, 1 ml Made in USA Lot No: 161898 Exp: 7/23/2013 Storage: <= -10 Degrees C Solvent: Methylene Chloride Motor oil composite Lot #: 161898 - 28615 Rec: 4/14/11 MFR exp. 07/23/13 OP: 9/1/11 EX: 9/1/12						
DIESEL 2ND SOURCE							
DIESEL FUEL #2	50,000mg/L	0281	1000mL	50mL	1000mg/L	MC # 051711B	9/1/11 EX: 3/1/12
	Diesel Fuel #2 Composite, 50,000 mg/L, 1 ml 911598-83 Lot # 167769 Storage 5-10 Degrees C Expiry 2/15/15 Made in the USA Solvent: Methylene Chloride Diesel Fuel #2 Composite OP: 9/1/11 Lot #: 167769 - 28397 EX: 9/1/12 Rec: 8/26/11 MFR exp. 02/15/15						

STANDARD | INITIAL CONC | SOURCE DATE | ALIQUOT | FINAL VOLUME | FINAL CONC | SOLVENT / LOT# | DATE / LOT#

KEROSENE/JP5 STD

STD	INITIAL CONC	SOURCE DATE	ALIQUOT	FINAL VOL.	FINAL CONC	SOLVENT / LOT#
JP5/ KEROSENE	50,000 µg/mL	O2SI CAT #010597-S50 LOT# 159381 OP:4/18/11 EX: 4/18/12	500 µL	25 mL	1000 µg/mL	MC LOT# 032811C

LAC 10/18/11 EX: 4/18/12

KEROSENE/JP5 CURVE

STANDARD	[µg/mL]	LOT #	DATE	EXP. DATE	µL	µL	µL	µL	µL	µL
Kerosene	1000		10/18/2011	4/18/2012	50	100	400	600	800	1000
JP5	MC	032811C			950	900	600	400	200	NA
				Final VOL.	1000	1000	1000	1000	1000	1000

LAC 10/18/11 EX: 4/18/12

DIESEL CCV 400ug/ml

STANDARD	INITIAL CONC.	SOURCE DATE	ALIQUOT	FINAL VOL	FINAL CONC.	SOLVENT /LOT#
DIESEL STD.	1000UG/ML	O2SI	400µL	1mL	400 µg/ml	MC
		9/1/2011	3/1/2012			051711B

LAC 10/18/11 EX: 3/1/12

MOTOR OIL CCV 400UG/ML

STANDARD	INITIAL CONC.	SOURCE DATE	ALIQUOT	FINAL VOL	FINAL CONC.	SOLVENT /LOT#
MOTOR OIL STD	1000UG/ML	O2SI	400µL	1mL	400 µg/ml	MC
		9/1/2011	3/1/2012			051711B

PRODIAMINE STANDARD

diamine | 1000µg/mL | O2SI | 10mL | 10mL | 1mL/mL | ACETONE

CAT: 031919-02
LOT: 161445-26939
OP: 10/19/11
EX: 7/1/12

#011011C

10/19/11 EX: 4/19/12

10/19/11

Prodiamine Curve

PREP DATE:	10/19/2011										
EXP:	4/19/2012										
SUPPLIER	ID#	µg/mL	LOT #	DATE	EXP. DATE	µL	µL	µL	µL	µL	µL
	Prodiamine	1		10/19/2011	4/19/2012	5	50	100	150	200	250
	HEXANE		082610B			995	950	900	850	800	750
					Final VOL.	1000	1000	1000	1000	1000	1000

10/19/11 EX: 4/19/12

STANDARD 082	INITIAL CONC	SOURCE DATE	ALIQUOT	FINAL VOLUME	FINAL CONC	SOLVENT LOT #	DATE / INITIALS

DIESEL SPIKE

DIESEL FUEL #2	50,000mg/L	O2SI	2000uL	50mL	2000mg/L	MC #51204	10/21/11 EX: 1/21/12

OP: 10/21/11
EX: 10/21/12
Diesel Fuel #2 Composite, 50,000 mg/L, 1 ml
011593-03
Lot# 179635 Storage ≤ 10 Degrees C Expiry 11/8/15
Diesel Fuel #2 Composite
Lot #: 179635 - 29641
Rec: 10/13/11 MFR exp. 11/08/15

OP: 10/21/11
EX: 10/21/12
Diesel Fuel #2 Composite, 50,000 mg/L, 1 ml
011593-03
Lot# 179635 Storage ≤ 10 Degrees C Expiry 11/8/15
Diesel Fuel #2 Composite
Lot #: 179635 - 29642
Rec: 10/13/11 MFR exp. 11/08/15

10/21/11

508 AMPULE * GAVE TO EXTRACTION TO SPIKE *

VARIOUS Analytes	100/2000mg/L	O2SI	N/A	1mL	100/2000mg/L	N/A	10/21/11 EX: 10/21/12

o2si
OCC Pesticide Standard, 100/2000 mg/L, 1 ml
Cat. No: 130200-02 Exp: 2/21/2013
Lot No: 156275 Storage: <= 6 Degrees C
Solvent: Tol.:Hex, 1:1
For Research Use Only
OCC Pesticide 100/2000mg/L
Lot #: 156275 - 26160
Rec: 2/23/10 MFR exp: 02/21/13
...ened: 10/21/11 EX: 10/21/12

10/21/11

		508 CALIBRATION CURVE				
Compound	Conc. In Mix	Conc. Of Stock	Aliquot	stock source	Final Vol.	Solvent Lot#
alachlor	(1) 0.005/0.1	5/100ug/ml	10ul	508 stock	10 mL	Hexane
benfluralin	(2) 0.03/0.6	5/100ug/ml	60ul	prep: 4/8/11	10 mL	# 082610B
captan	(3) 0.05/1.0	5/100ug/ml	250ul	Exp: 1/25/12	25 mL	
carbophenothion	(4) 0.1/2.0	5/100ug/ml	200ul		10 mL	
chlorothalonil	(5) 0.15/3.0	5/100ug/ml	300ul		10 mL	
chlorthal(dacthal)	(6) 0.2/4.0	5/100ug/ml	400ul		10 mL	
2,6 dichlorobenzonitrile(dicofol)						
kelthane						
nitrofen						
oxadiazon						
oxyfluorfen						
propachlor						
op DDD						
op DDE						
op DDT						
bis(2-ethylhexyl)phthalate						

10/21/11 EX: 1/25/12

10/21/11

		508 2ND SRC				
Compound	Init. Conc.	Stock Src	Aliquot	Final Vol	Final Conc.	Solvent Lot#
See Above	5/100 ug/ml	508 2nd Src Stock	250 uL	25 mL	0.05/1 ug/ml	Hexane
		Prep: 10/21/11				082610B
		Exp: 4/8/12				

10/21/11 EX: 4/8/12

STANDARD 088

INITIAL CONC	SOURCE	DATE	ALIQUOT	FINAL VOLUME	FINAL CONC	SOLVENT LOT #	DATE INITIALS

LAC 10/28/11

TCH SURROGATE CURVE

STD	[µg/mL]	LOT #	DATE	EXP. DATE	µL	µL	µL	µL	µL	µL
THC SURR	50	176405	10/17/2011	4/17/2012	50	100	400	600	800	1000
MC		61204			950	900	600	400	200	NA
				Final VOL.	1000	1000	1,000	1000	1000	1000

LAC 10/28/11 EX: 4/17/12

DIESEL CURVE

STD	[µg/mL]	LOT #	DATE	EXP. DATE	µL	µL	µL	µL	µL	µL
DIESEL	1000		10/26/2011	4/26/2012	10	100	400	600	800	1000
MC		51204			990	900	600	400	200	NA
				Final VOL.	1000	1000	1,000	1000	1000	1000

LAC 10/28/11 EX: 4/26/12

MOTOR OIL CURVE

STD	[µg/mL]	LOT #	DATE	EXP. DATE	µL	µL	µL	µL	µL	µL
MOTOR OIL	1000		10/26/2011	4/26/2012	50	100	400	600	800	1000
MC		51204			950	900	600	400	200	NA
				Final VOL.	1000	1000	1,000	1000	1000	1000

DIESEL 2ND SOURCE

STD	Init. Conc	Source	Aliquot	Final Vol.	Final Conc.	Solvent
DIESEL 2ND SRC	1000µg/ml	O2SI	400µL	1 mL	400 µg/mL	MC
	Prep:	9/1/2011				51204
	Exp:	3/1/2012				

LAC 10/28/11 EX: 3/1/12

LAC 10/28/11

KEROSENE/JP5 CURVE

STANDARD	[µg/mL]	LOT #	DATE	EXP. DATE	µL	µL	µL	µL	µL	µL
Kerosene	1000		10/18/2011	4/18/2012	50	100	400	600	800	1000
JP5	MC	51204			950	900	600	400	200	NA
				Final VOL.	1000	1000	1000	1000	1000	1000

LAC 10/28/11 EX: 4/18/12

11/1/11

PREP DATE:	11/1/2011					
OP 2ND SOURCE						
EXP:	4/19/2012					
SUPPLIER	ID#	[µg/mL]	LOT #	DATE	EXP. DATE	µL
	OP 2ND SRC	5		10/19/2011	4/19/2012	500
VWR	HEXANE		0826108			500
					Final VOL.	1000

11/1/11 EX: 4/19/12

PREP DATE:	11/1/2011											
OPF CURVE												
EXP:	4/13/2012											
						IA	1	2	3	4	5	6
SUPPLIER	ID#	[µg/mL]	LOT #	DATE	EXP. DATE	µL	µL	µL	µL	µL	µL	µL
	OPF STD	5		10/19/2011	4/13/2012	2	10	50	100	500	700	1000
	Hexane		0826108			998	990	950	900	500	300	NA
					Final VOL.	1000	1000	1000	1000	1000	1000	1000

11/1/11 EX: 4/13/12

PREP DATE:	11/1/2011										
OPC CURVE											
EXP:	3/15/2012										
						1	2	3	4	5	6
SUPPLIER	ID#	[µg/mL]	LOT #	DATE	EXP. DATE	µL	µL	µL	µL	µL	µL
	OPC STD	5		9/15/2011	3/15/2012	10	50	200	500	700	1000
	Hexane		0826108			990	950	800	500	300	NA
					Final VOL.	1000	1000	1000	1000	1000	1000

11/1/11

11/1/11 EX: 3/15/12

STANDARD	INITIAL CONC	SOURCE DATE	ALIQUOT	FINAL VOLUME	FINAL CONC	SOLVENT LOT #	DATE / INITIALS
		DIESEL STANDARD					
DIESEL FUEL #2	50,000ug/ml	O2SI	1000ul	50ML	1000ug/ml	MC # 51204	10/26/11 ex: 4/26/12
		Diesel Fuel #2 Composite, 50,000 mg/L, 1 ml; 111598-03; Lot # 167768; Storage 5-10 Degrees C; Expiry 2/15/15; Solvent Methylene Chloride; Made in the USA; Diesel Fuel #2 Composite OP: 10/26/11; Lot #: 167768 - 29406 EX: 10/26/12; Rec: 8/26/11 MFR exp. 02/15/15					
TERPHENYL HEXANE	600ug/ml	O2SI	4170ul		50ug/ml		
		CAT: 110316-05; LOT: 176405-29338; OP: 10/10/11; EX: 10/10/12					
		MOTOR OIL STANDARD					
MOTOR OIL	50,000ug/ml	O2SI	1000ul	50ML	1000ug/ml	MC # 51204	10/26/11 ex: 4/26/12
		OP: 10/26/11 EX: 10/26/12; o2si smart solutions; Motor Oil Composite, 50,000 mg/L, 1 ml; 116390-01; Storage: <= -10 Degrees C; Made in USA; Lot No: 161898; Solvent: Methylene Chloride; Exp: 7/23/2013; Motor oil composite; Lot #: 161898 - 28616; Rec: 4/14/11 MFR exp. 07/23/13					

		PAC ECO 2ND SOURCE			
DIAZINON	6ug/ml	200ug/ml	260ul	O2SI	10ml
DISULFOTON		200	CAT:	130169-01	HEXANE
MALATHION		200	LOT:	178204-28481	LOT#
MOLINATE		200	OP:	10/26/2011	082610B
PHORATE		200	EXP:	3/11/2012	
THIOBENCARB		200			
TRIBUTYL PHOSPHATE		200			
DEMETON		200			
DISCHLORVOS		200			
EPTC		200			
PARATHION		200			
AZINPHOS METHYL		200			
CHLORPYRIFOS		200			
DIMETHOATE		200			
METHIDATHION		200			
METHYL PARATHION		200			
ATRAZINE		200			
CYANIZINE		200			
TRIPHENYL PHOSPHATE		200			
PENDIMETHALIN (PROWL		200			
TRIFLURALIN		200			
SIMAZINE		200			

10/26/11
EX: 3/11/12

10/26/11

STANDARD

	INITIAL CONC	SOURCE DATE	ALIQUOT	FINAL VOLUME	FINAL CONC	SOLVENT LOT#	DATE: 0995

11/7/11

PREP:	11/7/2011												
PAC ECO CURVE													
EXP:	2/25/2012												
	ID#	[µg/mL]	LOT #	DATE	EXP. DATE	µL	µL	µL	µL	µL	µL	µL
	PAC ECO CAL STD	5		10/26/2011	2/25/2012	2	10	50	200	500	700	1000
VWR	Hexane		010711A			998	990	950	800	500	300	N/A
					Final VOL.	1000	1000	1000	1000	1000	1000	1000
PAC ECO 2ND SRC				DATE	EXP. DATE							
Prep: 11/7/11	Exp: 12/17/11	5	010711A	10/28/2011	12/17/2011	500/1000						

11/7/11 Ex: 2/25/12

11/7/11 Ex: 12/17/11

TCH SURROGATE CURVE

STD	[µg/mL]	LOT #	DATE	EXP. DATE	µL	µL	µL	µL	µL	µL
THC SURR	50	176405	10/17/2011	4/17/2012	50	100	400	600	800	1000
MC		51204			950	900	600	400	200	NA
				Final VOL.	1000	1000	1,000	1000	1000	1000

11/8/11 Ex: 4/17/12

DIESEL CURVE

STD	[µg/mL]	LOT #	DATE	EXP. DATE	µL	µL	µL	µL	µL	µL
DIESEL	1000		10/26/2011	4/26/2012	10	100	400	600	800	1000
MC		51204			990	900	600	400	200	NA
				Final VOL.	1000	1000	1,000	1000	1000	1000

11/8/11 Ex: 4/26/12

MOTOR OIL CURVE

STD	[µg/mL]	LOT #	DATE	EXP. DATE	µL	µL	µL	µL	µL	µL
MOTOR OIL	1000		10/26/2011	4/26/2012	50	100	400	600	800	1000
MC		51204			950	900	600	400	200	NA
				Final VOL.	1000	1000	1,000	1000	1000	1000

DIESEL 2ND SOURCE

STD	Init. Conc	Source	Aliquot	Final Vol.	Final Conc.	Solvent
DIESEL 2ND SRC	1000µg/ml	O2SI	400µL	1 mL	400 µg/mL	MC
	Prep:	9/1/2011				51204
	Exp:	3/1/2012				

11/8/11 Ex: 3/1/12

11/9/11

PREP DATE:	11/9/2011											
TERBACIL CURVE												
EXP:	3/13/2012											
						0.05	0.25	1	2.5	3.5	5	
SUPPLIER	ID#	[µg/mL]	LOT #	DATE	EXP. DATE	µL	µL	µL	µL	µL	µL	
	TERBACIL STD	5		9/13/2011	3/13/2012	10	50	200	500	700	1000	
VWR	HEXANE		082810B			950	920	900	800	500	300	
					Final VOL.	1000	1000	1000	1000	1000	1000	

11/9/11 Ex: 3/13/12

PREP DATE:	11/9/2011					
OP 2ND SOURCE						
EXP:	4/19/2012					
SUPPLIER	ID#	[µg/mL]	LOT #	DATE	EXP. DATE	µL
	OP 2ND SRC	5		10/19/2011	4/19/2012	500
VWR	HEXANE		082610B			500
					Final VOL.	1000

11/9/11 Ex: 4/19/12

11/9/11

PREP DATE:	11/9/2011											
OPF CURVE												
EXP:	2/7/2012											
						IA	1	2	3	4	5	6
SUPPLIER	ID#	[µg/mL]	LOT #	DATE	EXP. DATE	µL	µL	µL	µL	µL	µL	µL
	OPF STD	5		11/3/2011	2/7/2012	2	10	50	200	500	700	1000
	Hexane		082610B			998	990	950	800	500	300	NA
					Final VOL.	1000	1000	1000	1000	1000	1000	1000

11/9/11 Ex: 2/7/12

STANDARD	INITIAL CONC	SOURCE	DATE	ALIQUOT	FINAL VOLUME	FINAL CONC	SOLVENT / LOT #	DATE / INITIALS
				PCB SOIL SPIKE				
AR 1260	1000ng/mL	0251		1250mL	25mL	50ng/µL	ACETONE	
AR 1016	CAT: 130011-03						#	11/10/11
	LOT: 163607-27215							EX: 2/10/12
	OP: 11/10/11							
	EX: 11/10/12							
	AND							
	LOT: 152374-27210							
	OP: 3/2/11							
	EX: 3/2/12							
				PCB WATER SPIKE				
AR1016	1000ng/mL	0251		125mL	25mL	5ng/µL	ACETONE	
AR1260	CAT: 130011-03						#	11/10/11
	LOT: 163607-27214							EX: 2/10/12
	OP: 8/2/11							
	EX: 8/2/12							
				HERB 100/1000 (LVL 3) CCV				
VARIOUS	VARIOUS	HERB STD.		100µL	1mL	100ng/µL	MTBE	
SEE PG 075	PREP: 10/11/11						#	11/10/11
	EX: 4/11/12							EX: 4/11/12
				THC SURROGATE CAL. STD.				
D-TERPHENYL	1000ng/mL	0251		834µL	10mL	50ng/µL	MC	
DECACYCLENE	CAT: 110316-05						#	11/15/11
	LOT: 176405-29342							EX: 5/15/12
	OP: 10/10/11							
	EX: 10/10/12							

LAC 11/15/11

TCH SURROGATE CURVE											
STD	[µg/mL]	LOT #	DATE	EXP. DATE	µL	µL	µL	µL	µL	µL	LAC 11/15/11
THC SURR	50	176405	11/15/2011	5/15/2012	50	100	400	600	800	1000	EX 5/15/12
MC		51204			950	900	600	400	200	NA	
				Final VOL.	1000	1000	1,000	1000	1000	1000	

STANDARD | INITIAL CONC | SOURCE DATE | ALIQUOT | FINAL VOLUME | FINAL CONC | SOLVENT / LOT # | DATE / INITIALS

DIESEL CCV 400ug/ml							
STANDARD	INITIAL CONC.	SOURCE DATE	ALIQUOT	FINAL VOL	FINAL CONC.	SOLVENT /LOT#	
DIESEL STD.	1000UG/ML	O2SI	400µL	1mL	400 µg/ml	MC	LAC 11/29/11
		10/26/11	04/26/12			51204	EXP 4/26/11

MOTOR OIL CCV 400UG/ML						
STANDARD	INITIAL CONC.	SOURCE DATE	ALIQUOT	FINAL VOL	FINAL CONC.	SOLVENT /LOT#
MOTOR OIL STD	1000UG/ML	O2SI	400µL	1mL	400 µg/ml	MC
		10/26/11	04/26/12			51204

STANDARD	INITIAL CONC	SOURCE DATE	ALIQUOT	FINAL VOLUME	FINAL CONC	SOLVENT / LOT #	DATE / INITIALS

OCL/OP WATER SURROGATE

DECA — 3000 mg/L — O2SI — 30 mL — 100 mL — 1.5 mg/L — ACETONE # 011011C — 11/2/11 — EX: 2/2/12

TCMX

DBC — CAT: 130010-02 — LOT: 154164-29416 — OP: 11/2/11 — EX: 11/2/12

TBP — 1000 mg/L — O2SI — 500 mL — 5 mg/L

TPP

Tributyl- and Triphenylphosphate Solution, 1,000 mg/L, 1 ml
Cat# 130161-02
Lot# 164817 Storage 5-10 Degrees C Expiry 10/19/13
Solv: Acetone
Tributyl and Triphenyl phosphate
Lot #: 164817-27660
Rec: 10/20/10 MFR exp. 10/19/13

OP: 11/2/11 EX: 11/2/12 — 11/2/11

OCL/OP WATER SURROGATE

DECA — 5000 mg/L — O2SI — 30 mL — 100 mL — 1.5 mg/L — ACETONE # 011011C — 11/3/11 — EX: 2/3/12

TCMX

DBC — CAT: 130010-02 — LOT: 154164-29416 — OP: 11/2/11 — EX: 11/2/12

TBP — 1000 mg/L — O2SI — 500 mL — 5 mg/L

TPP — CAT: 130161-02 — LOT: 164817-27660 — OP: 11/2/11 — EX: 11/2/12

* WAS NOT RECORDED ON 11/2/11 — 11/3/11

DIESEL CCV 400ug/ml						
STANDARD	INITIAL CONC.	SOURCE DATE	ALIQUOT	FINAL VOL	FINAL CONC.	SOLVENT /LOT#
DIESEL STD.	1000UG/ML	O2SI	400µL	1mL	400 µg/ml	MC
		10/26/2011	4/26/2012			51204

11/2/11 — EX: 4/26/12

MOTOR OIL CCV 400UG/ML						
STANDARD	INITIAL CONC.	SOURCE DATE	ALIQUOT	FINAL VOL	FINAL CONC.	SOLVENT /LOT#
MOTOR OIL STD	1000UG/ML	O2SI	400µL	1mL	400 µg/ml	MC
		10/26/2011	4/26/2012			51204

KEROSENE CCV 400 UG/ML						
STANDARD	INITIAL CONC.	SOURCE DATE	ALIQUOT	FINAL VOL	FINAL CONC.	SOLVENT /LOT#
KEROSENE /JP5 STD	1000UG/ML	O2SI	400 µL	1ML	400 UG/ML	MC
		10/18/2011	4/18/2012		1mL	010611B

11/2/11 — EX: 4/18/12

Organic Extraction Worksheet

Method	THC-Separatory Funnel Extraction 3510C	Extraction Set	111031A	Extraction Method	SEP011	Units	mL

Spiked ID 1	Diesel Spike 10/21/11 EX 1/21/12		Surrogate ID 1	THC Surrogate 176405-29339
Spiked ID 2	Motor Oil Spike 8/5/11 EX 11/5/11		Surrogate ID 2	
Spiked ID 3			Surrogate ID 3	
Spiked ID 4			Surrogate ID 4	
Spiked ID 5			Surrogate ID 5	
Spiked ID 6			Sufficient Vol for Matrix QC:	YES
Spiked ID 7			Ext. Start Time:	
Spiked ID 8			Ext. End Time:	
			GC Requires Extract By:	11/02/11 0:00

pH1			Water Bath Temp Criteria	80 °C
pH2				
pH3				

Spiked By: HW **Date** 10/31/2011 **Witnessed By:** DL **Date** 10/31/2011

	Sample	Sample Container	Spike Amount	Spike ID	Surrogate Amount	Surrogate ID	Extract Amount	Final Volume	pH	Extract Date/Time	Comments
1	111031A Blk				0.250	1 equip	1000 E-WB5	5	7	10/31/11 15:00	
2	111031A LCS-1		1	1	0.250	1 equip	1000 E-WB5	5	7	10/31/11 15:00	
3	111031A LCS-2		1	2	0.250	1 equip	1000 E-WB5	5	7	10/31/11 15:00	
4	AY49327	AY49327W05			0.250	1 equip	1040 E-WB5	5	7	10/31/11 15:00	66103-1 WEEK RUSH -- Amber Liter
5	AY49328	AY49328W05			0.250	1 equip	1000 E-WB5	5	7	10/31/11 15:00	66103-1 WEEK RUSH -- Amber Liter
6	AY49329	AY49329W05			0.250	1 equip	1010 E-WB5	5	7	10/31/11 15:00	66103-1 WEEK RUSH -- Amber Liter
7	AY49330	AY49330W04			0.250	1 equip	1040 E-WB5	5	7	10/31/11 15:00	66103-1 WEEK RUSH -- Amber Liter
8	AY49331	AY49331W04			0.250	1 equip	1030 E-WB5	5	7	10/31/11 15:00	66103-1 WEEK RUSH -- Amber Liter
9	AY49333	AY49333W09			0.250	1 equip	1030 E-WB5	5	7	10/31/11 15:00	66102-2 WEEK RUSH -- Amber Liter
10	AY49334 MS-1	AY49334W35	1	1	0.250	1 equip	1030 E-WB6	5	7	10/31/11 15:00	66102-2 WEEK RUSH -- Amber Liter
11	AY49334 MSD-1	AY49334W36	1	1	0.250	1 equip	1030 E-WB6	5	7	10/31/11 15:00	66102-2 WEEK RUSH -- Amber Liter
12	AY49334	AY49334W32			0.250	1 equip	1030 E-WB5	5	7	10/31/11 15:00	66102-2 WEEK RUSH -- Amber Liter
13	AY49336	AY49336W09			0.250	1 equip	1050 E-WB6	5	7	10/31/11 15:00	66102-2 WEEK RUSH -- Amber Liter

Solvent and Lot#	
MC	BMD 51204
Na2SO4	3581C501

Extraction COC Transfer	
Extraction lab employee Initials	HW
GC analyst's initials	
Date	11/2/11
Time	11:30
Refrigerator	Hob met

	Technician's Initials
Scanned By	HW
Sample Preparation	HW
Extraction	HW/CC/DL
Concentration	JL

Modified	10/31/2011 2:12:11 PM

Reviewed By: HW 155 **Date** 11/1/2011

Organic Extraction Worksheet

Method	THC Separatory Funnel Extraction 3510C	Extraction Set	111031A	Extraction Method	SBP011	Units	mL

Spiked ID 1	Diesel Spike 10/21/11 EX 1/21/12		Surrogate ID 1	THC Surrogate 176405-29339	
Spiked ID 2	Motor Oil Spike 8/5/11 EX 11/5/11		Surrogate ID 2		
Spiked ID 3			Surrogate ID 3		
Spiked ID 4			Surrogate ID 4		
Spiked ID 5			Surrogate ID 5		
Spiked ID 6			Sufficient Vol for Matrix QC:	YES	
Spiked ID 7			Ext. Start Time:		
Spiked ID 8			Ext. End Time:		
			GC Requires Extract By:	11/02/11 0:00	

pH1			Water Bath Temp Criteria	80 °C
pH2				
pH3				

Spiked By: HW **Date** 10/31/2011 **Witnessed By:** DL **Date** 10/31/2011

	Sample	Sample Container	Spike Amount	Spike ID	Surrogate Amount	Surrogate ID	Extract Amount	Final Volume	pH	Extract Date/Time	Comments
14	AY49481	AY49481W10			0.250	1	1040	5	7	10/31/11 15:00	66116-2 WEEK RUSH -- Amber Liter
						equip	E-WB6				
15	AY49482	AY49482W10			0.250	1	1040	5	7	10/31/11 15:00	66116-2 WEEK RUSH -- Amber Liter
						equip	E-WB6				

HW 11/1/11

Solvent and Lot#	
MC	BMD 51204
Na2SO4	3581C501

Extraction COC Transfer	
Extraction lab employee Initials	HW
GC analyst's initials	
Date	11/2/11
Time	1120
Refrigerator	Upstairs

	Technician's Initials
Scanned By	HW
Sample Preparation	HW
Extraction	HW/CC/DL
Concentration	JL

Modified	10/31/2011 2:12:11 PM

Reviewed By: HW 156 **Date** 11/1/2011

Injection Log

Directory: G:\APOLLO\DATA\111028\111106\111108\111115\111129

Line	Vial	FileName	Multiplier	SampleName	Misc Info	Injected
1	3	1028003.D	1	DIESEL 10/1000 10/28/11	Mix(A)	10-28-11 9:47:18
2	4	1028004.D	1	DIESEL 100/1000	Mix(A)	10-28-11 10:11:19
3	5	1028005.D	1	DIESEL 400/1000	Mix(A)	10-28-11 10:35:26
4	6	1028006.D	1	DIESEL 600/1000	Mix(A)	10-28-11 10:59:35
5	7	1028007.D	1	DIESEL 800/1000	Mix(A)	10-28-11 11:23:49
6	8	1028008.D	1	DIESEL 1000/1000	Mix(A)	10-28-11 11:48:05
7	9	1028009.D	1	MOTOR OIL 50/1000 10/28/11	Mix(B)	10-28-11 12:12:27
8	10	1028010.D	1	MOTOR OIL 100/1000	Mix(B)	10-28-11 12:36:20
9	11	1028011.D	1	MOTOR OIL 400/1000	Mix(B)	10-28-11 13:00:16
10	12	1028012.D	1	MOTOR OIL 600/1000	Mix(B)	10-28-11 13:24:39
11	13	1028013.D	1	MOTOR OIL 800/1000	Mix(B)	10-28-11 13:48:43
12	14	1028014.D	1	MOTOR OIL 1000/1000	Mix(B)	10-28-11 14:13:14
13	15	1028015.D	1	DIESEL 2ND SRC 10/28/11	Mix(A)	10-28-11 14:37:14
14	16	1028016.D	1	THC SURR 10/1000 10/28/11	Mix(C)	10-28-11 15:01:44
15	17	1028017.D	1	THC SURR 100/1000	Mix(C)	10-28-11 15:25:58
16	18	1028018.D	1	THC SURR 400/1000	Mix(C)	10-28-11 15:50:20
17	19	1028019.D	1	THC SURR 600/1000	Mix(C)	10-28-11 16:14:52
18	20	1028020.D	1	THC SURR 800/1000	Mix(C)	10-28-11 16:38:57
19	21	1028021.D	1	THC SURR 1000/1000	Mix(C)	10-28-11 17:03:06
20	3	1106003.D	1	DIESEL 400/1000 10/28/11	Mix(A)	11-6-11 16:34:49
21	5	1106005.D	5	111031A BLK 5/1000	Water	11-6-11 17:22:09
22	13	1106013.D	4.85437	AY49333W09 5/1030	Water	11-6-11 20:30:17
23	14	1106014.D	4.85437	AY49334W35 MS-1 5/1030	Water	11-6-11 20:53:41
24	15	1106015.D	4.85437	AY49334W36 MSD-1 5/1030	Water	11-6-11 21:17:05
25	16	1106016.D	4.85437	AY49334W32 5/1030	Water	11-6-11 21:40:27
26	17	1106017.D	1	DIESEL 400/1000 11/2/11	Mix(A)	11-6-11 22:03:47
27	19	1106019.D	4.7619	AY49336W09 5/1050	Water	11-6-11 22:50:22
28	27	1106027.D	1	DIESEL 400/1000 11/2/11	Mix(A)	11-7-11 1:56:08
29	5	1108005.D	1	DIESEL 100/1000	Mix(A)	11-8-11 15:50:59
30	6	1108006.D	1	DIESEL 400/1000	Mix(A)	11-8-11 16:14:36
31	7	1108007.D	1	DIESEL 600/1000	Mix(A)	11-8-11 16:38:14
32	8	1108008.D	1	DIESEL 800/1000	Mix(A)	11-8-11 17:01:53
33	9	1108009.D	1	DIESEL 1000/1000	Mix(A)	11-8-11 17:25:32
34	11	1108011.D	1	MOTOR OIL 50/1000 11/8/11	Mix(B)	11-8-11 18:12:45
35	12	1108012.D	1	MOTOR OIL 100/1000	Mix(B)	11-8-11 18:36:14
36	13	1108013.D	1	MOTOR OIL 400/1000	Mix(B)	11-8-11 18:59:47
37	14	1108014.D	1	MOTOR OIL 600/1000	Mix(B)	11-8-11 19:23:20
38	15	1108015.D	1	MOTOR OIL 800/1000	Mix(B)	11-8-11 19:46:53
39	16	1108016.D	1	MOTOR OIL 1000/1000	Mix(B)	11-8-11 20:10:21
40	69	1108069.D	1	DIESEL 10/1000 11/8/11	Mix(A)	11-9-11 17:18:58
41	70	1108070.D	1	DIESEL 400 2ND SRC 11/8/11	Mix(A)	11-9-11 17:42:38
42	21	1115021.D	1	THC SURR 10/1000 11/15/11	Mix(C)	11-15-11 18:21:35
43	22	1115022.D	1	THC SURR 100/1000	Mix(C)	11-15-11 18:45:31
44	23	1115023.D	1	THC SURR 400/1000	Mix(C)	11-15-11 19:09:25
45	24	1115024.D	1	THC SURR 600/1000	Mix(C)	11-15-11 19:33:17
46	25	1115025.D	1	THC SURR 800/1000	Mix(C)	11-15-11 19:57:06
47	26	1115026.D	1	THC SURR 1000/1000	Mix(C)	11-15-11 20:20:52
48	12	1129012.D	1	DIESEL 400/1000 11/29/11	Mix(A)	11-29-11 13:06:20
49	17	1129017.D	5	111031A LCS-1 5/1000	Water	11-29-11 18:45:15
50	24	1129024.D	1	DIESEL 400/1000 11/29/11	Mix(A)	11-29-11 21:29:15

EPA METHOD 8270
Polynuclear Aromatic Hydrocarbons

APPL, INC.

EPA METHOD 8270
Polynuclear Aromatic Hydrocarbons
QC Summary

APPL, INC.

Method Blank
EPA 8270D SIM

Blank Name/QCG: **111031W-49334 - 161019**
Batch ID: #SIMHC-111031A

APPL Inc.
908 North Temperance Avenu
Clovis, CA 93611

Sample Type	Analyte	Result	LOQ	LOD	DL	Units	Extraction Date	Analysis Date
BLANK	1-METHYLNAPHTHALENE	0.12 U	0.2	0.12	0.06	ug/L	10/31/11	11/05/11
BLANK	2-METHYLNAPHTHALENE	0.12 U	0.2	0.12	0.06	ug/L	10/31/11	11/05/11
BLANK	ACENAPHTHENE	0.12 U	0.2	0.12	0.06	ug/L	10/31/11	11/05/11
BLANK	ACENAPHTHYLENE	0.12 U	0.2	0.12	0.06	ug/L	10/31/11	11/05/11
BLANK	ANTHRACENE	0.10 U	0.2	0.10	0.05	ug/L	10/31/11	11/05/11
BLANK	BENZO(A)ANTHRACENE	0.14 U	0.2	0.14	0.07	ug/L	10/31/11	11/05/11
BLANK	BENZO(A)PYRENE	0.14 U	0.2	0.14	0.07	ug/L	10/31/11	11/05/11
BLANK	BENZO(B)FLUORANTHENE	0.12 U	0.2	0.12	0.06	ug/L	10/31/11	11/05/11
BLANK	BENZO(GHI)PERYLENE	0.16 U	0.2	0.16	0.08	ug/L	10/31/11	11/05/11
BLANK	BENZO(K)FLUORANTHENE	0.14 U	0.2	0.14	0.07	ug/L	10/31/11	11/05/11
BLANK	CHRYSENE	0.10 U	0.2	0.10	0.05	ug/L	10/31/11	11/05/11
BLANK	DIBENZ(A,H)ANTHRACENE	0.10 U	0.2	0.10	0.05	ug/L	10/31/11	11/05/11
BLANK	FLUORANTHENE	0.16 U	0.2	0.16	0.08	ug/L	10/31/11	11/05/11
BLANK	FLUORENE	0.12 U	0.2	0.12	0.06	ug/L	10/31/11	11/05/11
BLANK	INDENO(1,2,3-CD)PYRENE	0.14 U	0.2	0.14	0.07	ug/L	10/31/11	11/05/11
BLANK	NAPHTHALENE	0.10 U	0.2	0.10	0.05	ug/L	10/31/11	11/05/11
BLANK	PHENANTHRENE	0.14 U	0.2	0.14	0.07	ug/L	10/31/11	11/05/11
BLANK	PYRENE	0.16 U	0.2	0.16	0.08	ug/L	10/31/11	11/05/11
BLANK	SURROGATE: 2-FLUORBIPHENY	51.7	50-110			%	10/31/11	11/05/11
BLANK	SURROGATE: NITROBENZENE-	66.3	40-110			%	10/31/11	11/05/11
BLANK	SURROGATE: TERPHENYL-D14 (	54.5	50-135			%	10/31/11	11/05/11

Quant Method:SIM2.M
Run #:1105L028
Instrument:Linus
Sequence:L111027
Initials:LF

GC SC-Blank-REG MDLs
Printed: 11/09/11 3:58:52 PM

Form 2 & 8

Surrogate Recovery

Lab Name: APPL, Inc. SDG No: 66102

Case No: 66102 Date Analyzed: 11/05/11

Matrix: WATER Instrument: Linus

APPL ID.	Client Sample No.	SURROGATE: 2-FLUORBIPHENYL (S)			SURROGATE: NITROBENZENE-D5 (S)		
		Limits	Result	Qualifier	Limits	Result	Qualifier
111031A-BLK	Blank	50-110	51.7		40-110	66.3	
111031A-LCS	Lab Control Spike	50-110	55.5		40-110	57.0	
AY49333	ES046	50-110	54.0		40-110	61.9	
AY49334-MS	Matrix Spike	50-110	53.6		40-110	59.3	
AY49334-MSD	Matrix SpikeD	50-110	55.7		40-110	53.1	
AY49334	ES047	50-110	53.4		40-110	65.4	
AY49336	ES049	50-110	65.0		40-110	64.4	

Comments: Batch: #SIMHC-111031A

Form 2 & 8

Surrogate Recovery

Lab Name: APPL, Inc. SDG No: 66102

Case No: 66102 Date Analyzed: 11/05/11

Matrix: WATER Instrument: Linus

APPL ID.	Client Sample No.	SURROGATE: TERPHENYL-D14 (S)					
		Limits	Result	Qualifier	Limits	Result	Qualifier
111031A-BLK	Blank	50-135	54.5				
111031A-LCS	Lab Control Spike	50-135	53.0				
AY49333	ES046	50-135	50.3				
AY49334-MS	Matrix Spike	50-135	50.7				
AY49334-MSD	Matrix SpikeD	50-135	61.9				
AY49334	ES047	50-135	61.4				
AY49336	ES049	50-135	51.9				

Comments: Batch: #SIMHC-111031A

Laboratory Control Spike Recovery

EPA 8270D SIM

APPL ID: **111031W-49334 LCS - 161019**
Batch ID: #SIMHC-111031A

APPL Inc.
908 North Temperance Avenue
Clovis, CA 93611

Compound Name	Spike Level ug/L	SPK Result ug/L	SPK % Recovery	Recovery Limits
1-METHYLNAPHTHALENE	4.00	2.25	56.3	45-105
2-METHYLNAPHTHALENE	4.00	2.26	56.5	45-105
ACENAPHTHENE	4.00	2.55	63.7	45-110
ACENAPHTHYLENE	4.00	2.39	59.8	50-105
ANTHRACENE	4.00	2.47	61.8	55-110
BENZO(A)ANTHRACENE	4.00	2.74	68.5	55-110
BENZO(A)PYRENE	4.00	2.48	62.0	55-110
BENZO(B)FLUORANTHENE	4.00	2.43	60.8	45-120
BENZO(GHI)PERYLENE	4.00	2.80	70.0	40-125
BENZO(K)FLUORANTHENE	4.00	3.23	80.8	45-125
CHRYSENE	4.00	2.86	71.5	55-110
DIBENZ(A,H)ANTHRACENE	4.00	2.89	72.3	40-125
FLUORANTHENE	4.00	2.86	71.5	55-115
FLUORENE	4.00	2.59	64.8	50-110
INDENO(1,2,3-CD)PYRENE	4.00	2.95	73.8	45-125
NAPHTHALENE	4.00	2.30	57.5	40-100
PHENANTHRENE	4.00	2.43	60.8	50-115
PYRENE	4.00	2.57	64.3	50-130
SURROGATE: 2-FLUORBIPHENYL (S)	2.00	1.11	55.5	50-110
SURROGATE: NITROBENZENE-D5 (S)	2.00	1.14	57.0	40-110
SURROGATE: TERPHENYL-D14 (S)	2.00	1.06	53.0	50-135

Comments: _____

Primary	SPK
Quant Method :	SIM2.M
Extraction Date :	10/31/11
Analysis Date :	11/05/11
Instrument :	Linus
Run :	1105L029
Initials :	LF

Printed: 11/09/11 3:59:06 PM

APPL Standard LCS

Matrix Spike Recoveries

EPA 8270D SIM

APPL ID: **111031W-49334 MS - 161019**
Batch ID: #SIMHC-111031A
Sample ID: AY49334
Client ID: ES047

APPL Inc.
908 North Temperance Avenue
Clovis, CA 93611

Compound Name	Spike Lvl ug/L	Matrix Result ug/L	SPK Result ug/L	DUP Result ug/L	SPK % Recovery	DUP % Recovery	Recovery Limits	RPD %	RPD Limits
1-METHYLNAPHTHALENE	3.88	0.62	2.88	2.71	58.2	53.9	45-105	6.1	25
2-METHYLNAPHTHALENE	3.88	0.20	2.26	1.98	53.1	45.9	45-105	13.2	25
ACENAPHTHENE	3.88	0.16	2.17	1.79	51.8	42.0 #	45-110	19.2	25
ACENAPHTHYLENE	3.88	ND	2.14	1.77	55.2	45.6 #	50-105	18.9	25
ANTHRACENE	3.88	ND	2.14	1.96	55.2	50.5 #	55-110	8.8	25
BENZO(A)ANTHRACENE	3.88	ND	3.04	2.84	78.4	73.2	55-110	6.8	25
BENZO(A)PYRENE	3.88	ND	2.62	2.30	67.5	59.3	55-110	13.0	25
BENZO(B)FLUORANTHENE	3.88	ND	2.72	2.34	70.1	60.3	45-120	15.0	25
BENZO(GHI)PERYLENE	3.88	ND	2.84	2.54	73.2	65.5	40-125	11.2	25
BENZO(K)FLUORANTHENE	3.88	ND	3.07	2.84	79.1	73.2	45-125	7.8	25
CHRYSENE	3.88	ND	2.75	2.56	70.9	66.0	55-110	7.2	25
DIBENZ(A,H)ANTHRACENE	3.88	ND	2.90	2.51	74.7	64.7	40-125	14.4	25
FLUORANTHENE	3.88	ND	2.84	2.59	73.2	66.8	55-115	9.2	25
FLUORENE	3.88	0.083	2.42	2.08	60.2	51.5	50-110	15.1	25
INDENO(1,2,3-CD)PYRENE	3.88	ND	2.93	2.71	75.5	69.8	45-125	7.8	25
NAPHTHALENE	3.88	1.0	3.70	2.56	69.6	40.2	40-100	36.4 #	25
PHENANTHRENE	3.88	ND	2.38	2.20	61.3	56.7	50-115	7.9	25
PYRENE	3.88	ND	2.63	2.45	67.8	63.1	50-130	7.1	25
SURROGATE: 2-FLUORBIPHENYL (S)	1.94	NA	1.04	1.08	53.6	55.7	50-110		
SURROGATE: NITROBENZENE-D5 (S)	1.94	NA	1.15	1.03	59.3	53.1	40-110		
SURROGATE: TERPHENYL-D14 (S)	1.94	NA	0.984	1.20	50.7	61.9	50-135		

= Recovery is outside QC limits.

Comments:

Primary	SPK	DUP
Quant Method :	SIM2.M	SIM2.M
Extraction Date :	10/31/11	10/31/11
Analysis Date :	11/05/11	11/06/11
Instrument :	Linus	Linus
Run :	1105L036	1105L037
Initials :	LF	

Printed: 11/09/11 4:01:33 PM
APPL MSD SCII

8270D-SIM

Form 4

Blank Summary

Lab Name: APPL, Inc. SDG No: 66102

Case No: 66102 Date Analyzed: 11/05/11

Matrix: WATER Instrument: Linus

Blank ID: 111031A-BLK Time Analyzed: 2015

APPL ID.	Client Sample No.	File ID.	Date Analyzed
111031A-BLK	Blank	1105L028	11/05/11 2015
111031A-LCS	Lab Control Spike	1105L029	11/05/11 2041
AY49333	ES046	1105L035	11/05/11 2311
111031A-MS	Matrix Spike	1105L036	11/05/11 2336
111031A-MSD	Matrix SpikeD	1105L037	11/06/11 0001
AY49334	ES047	1105L038	11/06/11 0026
AY49336	ES049	1105L039	11/06/11 0051

Comments: Batch: #SIMHC-111031A

Form 5
Tune Summary

Lab Name: APPL Inc.
Case No: 66102
Matrix: Water
ID: SVTUNE 10-27-11

SDG No: 66102
Date Analyzed: 11/05/11
Instrument: Linus
Time Analyzed: 16:36

	Client Sample No.	APPL ID.	File ID.	Date Analyzed
1	Blank	111031A BLK 1/1000	1105L028.D	11/05/11 20:15
2	Lab Control Spike	111031A LCS-1 1/1000	1105L029.D	11/05/11 20:41
3	ES046	AY49333W10 1/1050	1105L035.D	11/05/11 23:11
4	Matrix Spike	AY49334W30 MS-1 1/10	1105L036.D	11/05/11 23:36
5	Matrix Spike Dup	AY49334W34 MSD-1 1/1	1105L037.D	11/06/11 0:01
6	ES047	AY49334W29 1/1050	1105L038.D	11/06/11 0:26
7	ES049	AY49336W10 1/1030	1105L039.D	11/06/11 0:51
8				
9				
10				
11				
12				
13				
14				
15				
16				
17				
18				
19				
20				
21				
22				

m/e		
51	29.95 - 60% of mass 198	59.0
68	0 - 2.05% of mass 69	0.0
70	0 - 2% of mass 69	0.6
127	40 - 60% of mass 198	55.1
197	0 - 1% of mass 198	0.5
198	100 - 100% of mass 198	100.0
199	5 - 9% of mass 198	7.1
275	10 - 30% of mass 198	23.2
365	1 - 100% of mass 198	1.9
441	0.01 - 100% of mass 443	73.0
442	40 - 150% of mass 198	58.8
443	17 - 23% of mass 442	19.7

8A
INTERNAL STANDARD AREA AND RT SUMMARY

Lab Name: APPL Inc. Contract: Review

Lab Code: SDG No.: 66102

Lab File ID (Standard): 1028L007.D Date Analyzed: 10/28/11

Instrument ID: Linus Time Analyzed: 11:58

GC Column: ID: Heated Purge: (Y/N)

		Napthalene-D8(IS)		Acenaphthene-D10(IS)		Phenanthrene-D10(IS)	
		AREA #	RT #	AREA #	RT #	AREA #	RT #
	12 HOUR STD	2479	6.12	1083	8.11	1851	9.85
	UPPER LIMIT	4958	6.62	2166	8.61	3702	10.35
	LOWER LIMIT	1240	5.62	542	7.61	926	9.35
	SAMPLE NO.						
01	111031A BLK 1/1000	2305	6.12	1068	8.11	2122	9.86
02	111031A LCS-1 1/1000	2079	6.12	961	8.11	1713	9.85
03	AY49333W10 1/1050	2476	6.11	1226	8.11	1995	9.83
04	AY49334W30 MS-1 1/1(	2236	6.11	1033	8.11	1791	9.84
05	AY49334W34 MSD-1 1/	2287	6.11	1131	8.11	1879	9.83
06	AY49334W29 1/1050	2328	6.11	1134	8.11	1938	9.84
07	AY49336W10 1/1030	2268	6.12	1052	8.11	1894	9.86
08							
09							
10							
11							
12							
13							
14							
15							
16							
17							
18							
19							
20							
21							
22							

AREA UPPER LIMIT = +100% of internal standard area.
AREA LOWER LIMIT = -50% of internal standard area.
RT UPPER LIMIT = +0.50 minutes of internal standard RT
RT LOWER LIMIT = -0.50 minutes of internal standard RT

Column used to flag values outside QC limits with an asterisk.
* Values outside of QC limits.

8A
INTERNAL STANDARD AREA AND RT SUMMARY

Lab Name: APPL Inc. Contract: Review

Lab Code: _____ SDG No.: 66102

Lab File ID (Standard): 1028L007.D Date Analyzed: 10/28/11

Instrument ID: Linus Time Analyzed: 11:58

GC Column: _____ ID: Heated Purge: (Y/N) _____

		Chrysene-D12(IS)		Perylene-D12(IS)			
		AREA #	RT #	AREA #	RT #	AREA #	RT #
	12 HOUR STD	2378	12.93	1871	14.56		
	UPPER LIMIT	4756	13.43	3742	15.06		
	LOWER LIMIT	1189	12.43	936	14.06		
	SAMPLE NO.						
01	111031A BLK 1/1000	2454	12.94	2143	14.57		
02	111031A LCS-1 1/1000	2367	12.93	2017	14.56		
03	AY49333W10 1/1050	2680	12.94	2235	14.57		
04	AY49334W30 MS-1 1/1(	2429	12.93	2046	14.56		
05	AY49334W34 MSD-1 1/	2428	12.93	2160	14.56		
06	AY49334W29 1/1050	2444	12.94	2139	14.57		
07	AY49336W10 1/1030	2526	12.94	2131	14.57		
08							
09							
10							
11							
12							
13							
14							
15							
16							
17							
18							
19							
20							
21							
22							

AREA UPPER LIMIT = +100% of internal standard area.
AREA LOWER LIMIT = -50% of internal standard area.
RT UPPER LIMIT = +0.50 minutes of internal standard RT
RT LOWER LIMIT = -0.50 minutes of internal standard RT

Column used to flag values outside QC limits with an asterisk.
* Values outside of QC limits.

EPA METHOD 8270
Polynuclear Aromatic Hydrocarbons
Sample Data

APPL, INC.

EPA 8270D SIM

Environet, Inc.
650 Iwilei Rd, #204
Honolulu, HI 96817

APPL Inc.
908 North Temperance Avenue
Clovis, CA 93611

Attn: Stacy Fineran
Project: RED HILL/1022-024
Sample ID: ES046
Sample Collection Date: 10/24/11

ARF: 66102
APPL ID: AY49333
QCG: #SIMHC-111031A-161019

Method	Analyte	Result	LOQ	LOD	DL	Units	Extraction Date	Analysis Date
8270D-SIM	1-METHYLNAPHTHALENE	0.53	0.2	0.12	0.06	ug/L	10/31/11	11/05/11
8270D-SIM	2-METHYLNAPHTHALENE	0.15 J	0.2	0.12	0.06	ug/L	10/31/11	11/05/11
8270D-SIM	ACENAPHTHENE	0.17 J	0.2	0.12	0.06	ug/L	10/31/11	11/05/11
8270D-SIM	ACENAPHTHYLENE	0.12 U	0.2	0.12	0.06	ug/L	10/31/11	11/05/11
8270D-SIM	ANTHRACENE	0.10 U	0.2	0.10	0.05	ug/L	10/31/11	11/05/11
8270D-SIM	BENZO(A)ANTHRACENE	0.14 U	0.2	0.14	0.07	ug/L	10/31/11	11/05/11
8270D-SIM	BENZO(A)PYRENE	0.14 U	0.2	0.14	0.07	ug/L	10/31/11	11/05/11
8270D-SIM	BENZO(B)FLUORANTHENE	0.12 U	0.2	0.12	0.06	ug/L	10/31/11	11/05/11
8270D-SIM	BENZO(GHI)PERYLENE	0.16 U	0.2	0.16	0.08	ug/L	10/31/11	11/05/11
8270D-SIM	BENZO(K)FLUORANTHENE	0.14 U	0.2	0.14	0.07	ug/L	10/31/11	11/05/11
8270D-SIM	CHRYSENE	0.10 U	0.2	0.10	0.05	ug/L	10/31/11	11/05/11
8270D-SIM	DIBENZ(A,H)ANTHRACENE	0.10 U	0.2	0.10	0.05	ug/L	10/31/11	11/05/11
8270D-SIM	FLUORANTHENE	0.16 U	0.2	0.16	0.08	ug/L	10/31/11	11/05/11
8270D-SIM	FLUORENE	0.082 J	0.2	0.12	0.06	ug/L	10/31/11	11/05/11
8270D-SIM	INDENO(1,2,3-CD)PYRENE	0.14 U	0.2	0.14	0.07	ug/L	10/31/11	11/05/11
8270D-SIM	NAPHTHALENE	0.80	0.2	0.10	0.05	ug/L	10/31/11	11/05/11
8270D-SIM	PHENANTHRENE	0.14 U	0.2	0.14	0.07	ug/L	10/31/11	11/05/11
8270D-SIM	PYRENE	0.16 U	0.2	0.16	0.08	ug/L	10/31/11	11/05/11
8270D-SIM	SURROGATE: 2-FLUORBIPHENYL (S)	54.0	50-110			%	10/31/11	11/05/11
8270D-SIM	SURROGATE: NITROBENZENE-D5 (S)	61.9	40-110			%	10/31/11	11/05/11
8270D-SIM	SURROGATE: TERPHENYL-D14 (S)	50.3	50-135			%	10/31/11	11/05/11

J = Estimated value.

Quant Method: SIM2.M
Run #: 1105L035
Instrument: Linus
Sequence: L111027
Dilution Factor: 1
Initials: LF

Printed: 11/09/11 3:59:14 PM
APPL-F1-SC-NoMC-REG MDLs

Quantitation Report (QT Reviewed)

Data File : M:\LINUS\DATA\L111027\1105L035.D
Acq On : 5 Nov 11 23:11
Sample : AY49333W10 1/1050
Misc :

Vial: 35
Operator: LF
Inst : Linus
Multiplr: 0.95

Quant Time: Nov 9 9:17 2011 Quant Results File: SIM2.RES

Quant Method : M:\LINUS\DATA\L111027\SIM2.M (RTE Integrator)
Title : EPA 8270C
Last Update : Wed Nov 02 15:56:51 2011
Response via : Initial Calibration
DataAcq Meth : 87SIMAQ

Internal Standards	R.T.	QIon	Response	Conc Units	Dev(Min)
1) Napthalene-D8(IS)	6.11	136	2476	2.50000 ppb	-0.01
6) Acenaphthene-D10(IS)	8.11	164	1226	2.50000 ppb	0.00
11) Phenanthrene-D10(IS)	9.83	188	1995	2.50000 ppb	-0.01
15) Chrysene-D12(IS)	12.94	240	2680	2.50000 ppb	0.01
21) Perylene-D12(IS)	14.57	264	2235	2.50000 ppb	0.01
System Monitoring Compounds					
2) Surrogate Recovery (NBZ)	5.41	82	549	1.17909 ppb	-0.04
Spiked Amount 1.905			Recovery = 61.898%		
7) Surrogate Recovery (FBP)	7.35	172	1180	1.02825 ppb	0.00
Spiked Amount 1.905			Recovery = 53.970%		
17) Surrogate Recovery (TPH)	11.70	244	1161	0.95815 ppb	-0.01
Spiked Amount 1.905			Recovery = 50.295%		
Target Compounds					Qvalue
3) Naphthalene	6.13	128	1456	0.80384 ppb #	38
4) 2-Methylnaphthalene	6.93	142	142	0.15289 ppb	99
5) 1-Methylnaphthalene	7.02	142	568	0.52981 ppb	94
9) Acenaphthene	8.15	154	168	0.17136 ppb	92
10) Fluorene	8.75	166	88	0.08203 ppb	92

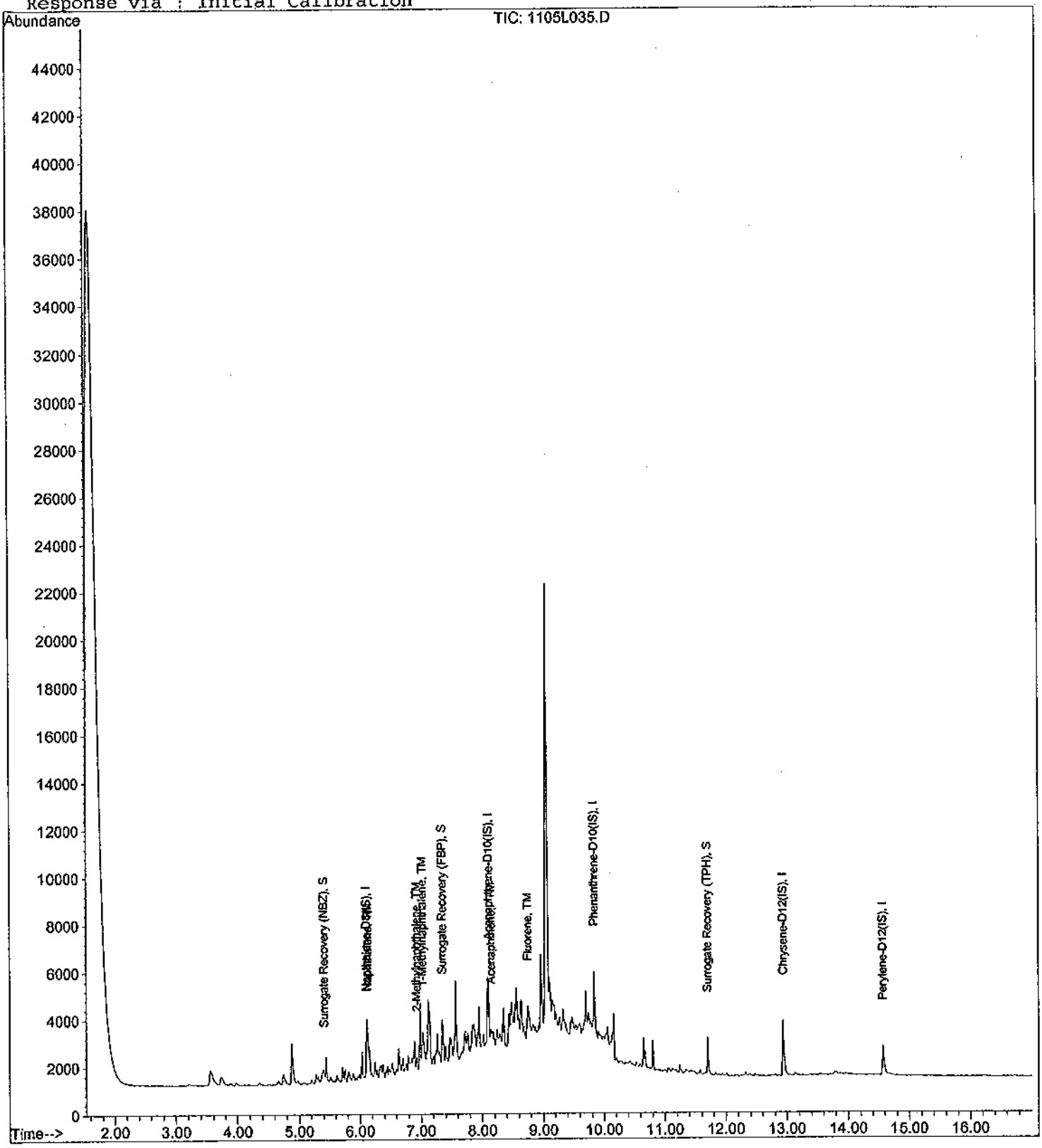
(#) = qualifier out of range (m) = manual integration

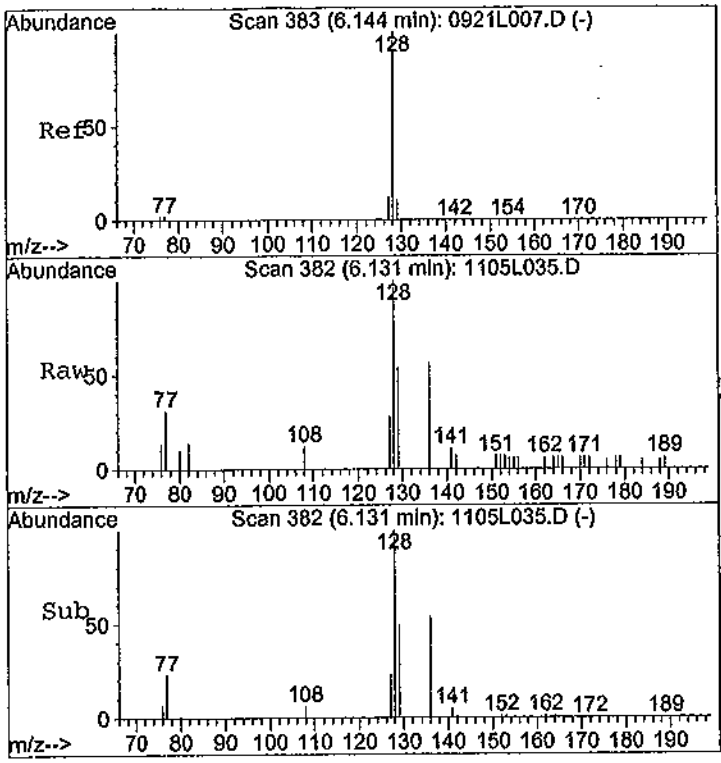
Quantitation Report

Data File : M:\LINUS\DATA\L111027\1105L035.D Vial: 35
Acq On : 5 Nov 11 23:11 Operator: LF
Sample : AY49333W10 1/1050 Inst : Linus
Misc : Multiplr: 0.95

Quant Time: Nov 9 9:17 2011 Quant Results File: SIM2.RES

Method : M:\LINUS\DATA\L111027\SIM2.M (RTE Integrator)
Title : EPA 8270C
Last Update : Tue Nov 08 16:22:04 2011
Response via : Initial Calibration

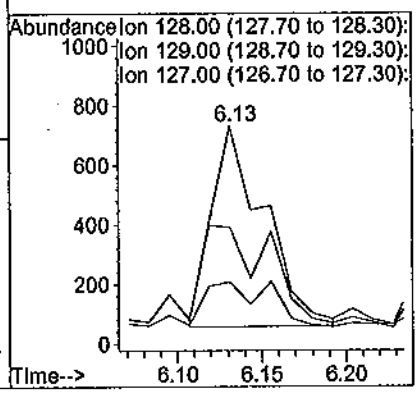


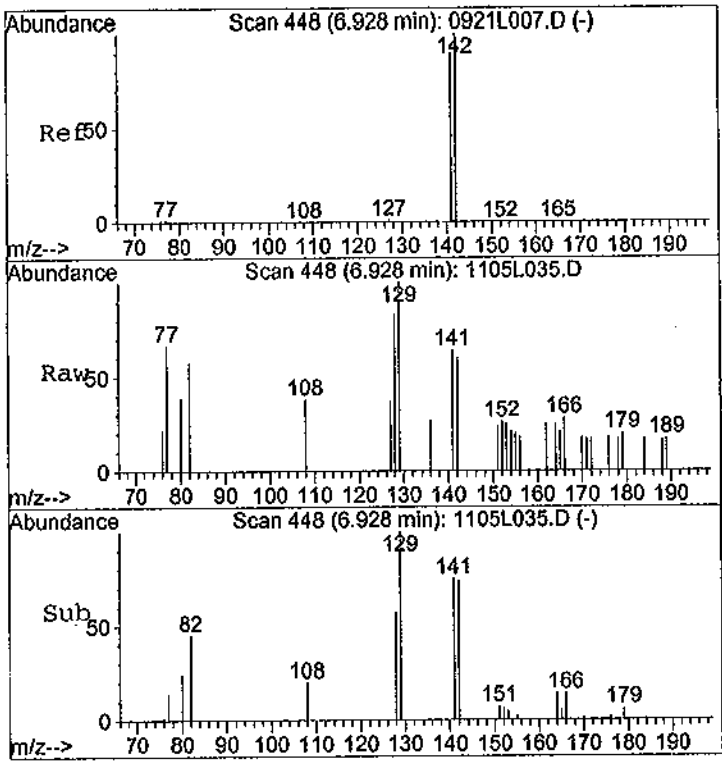


#3
Naphthalene
Concen: 0.80384 ppb
RT: 6.13 min Scan# 382
Delta R.T. -0.01 min
Lab File: 1105L035.D
Acq: 5 Nov 11 23:11

Tgt Ion:128 Resp: 1456

Ion	Ratio	Lower	Upper
128	100		
129	49.4	7.7	14.3#
127	23.0	8.6	16.0#

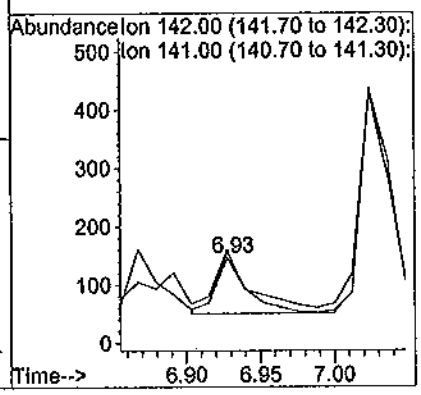


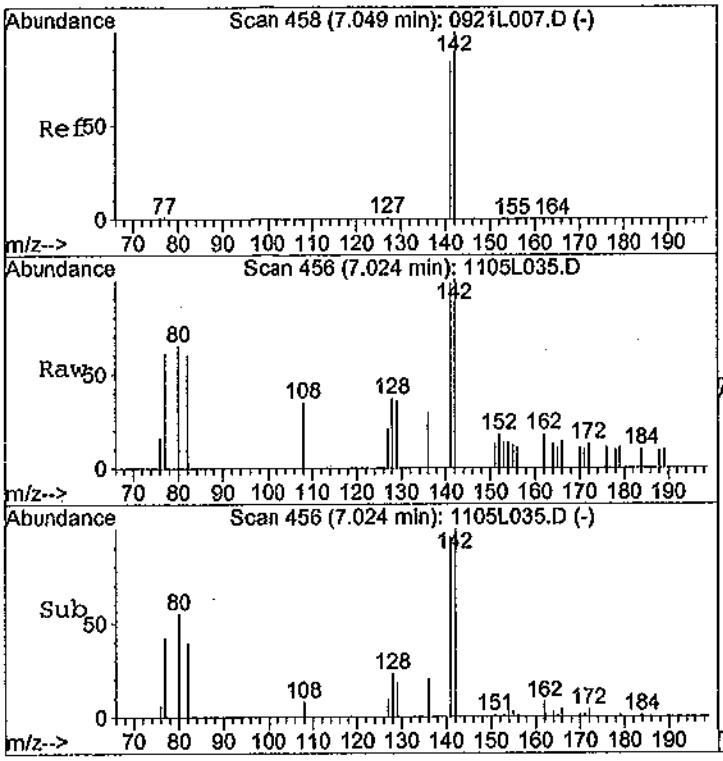


#4
2-Methylnaphthalene
Concen: 0.15289 ppb
RT: 6.93 min Scan# 448
Delta R.T. -0.04 min
Lab File: 1105L035.D
Acq: 5 Nov 11 23:11

Tgt Ion:142 Resp: 142

Ion	Ratio	Lower	Upper
142	100		
141	98.9	68.9	127.9

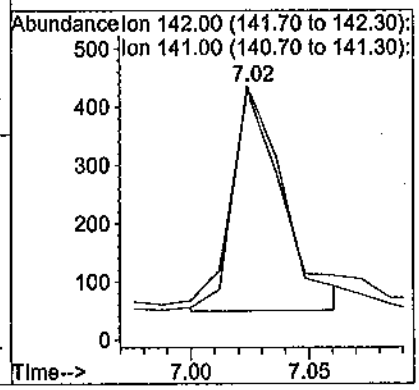


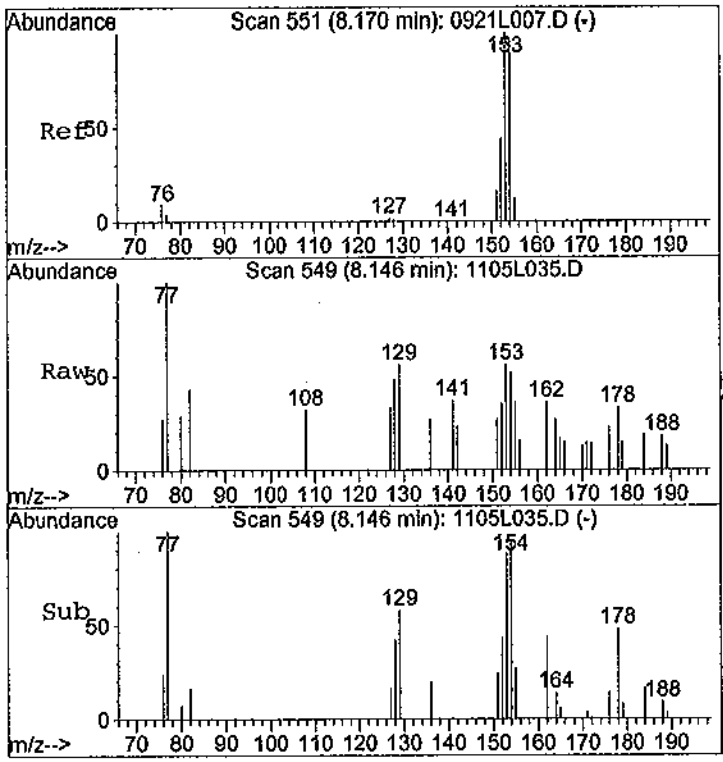


#5
1-Methylnaphthalene
Concen: 0.52981 ppb
RT: 7.02 min Scan# 456
Delta R.T. -0.02 min
Lab File: 1105L035.D
Acq: 5 Nov 11 23:11

Tgt Ion:142 Resp: 568

Ion	Ratio	Lower	Upper
142	100		
141	94.5	70.0	130.0

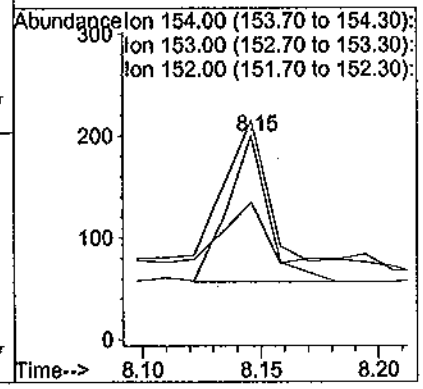


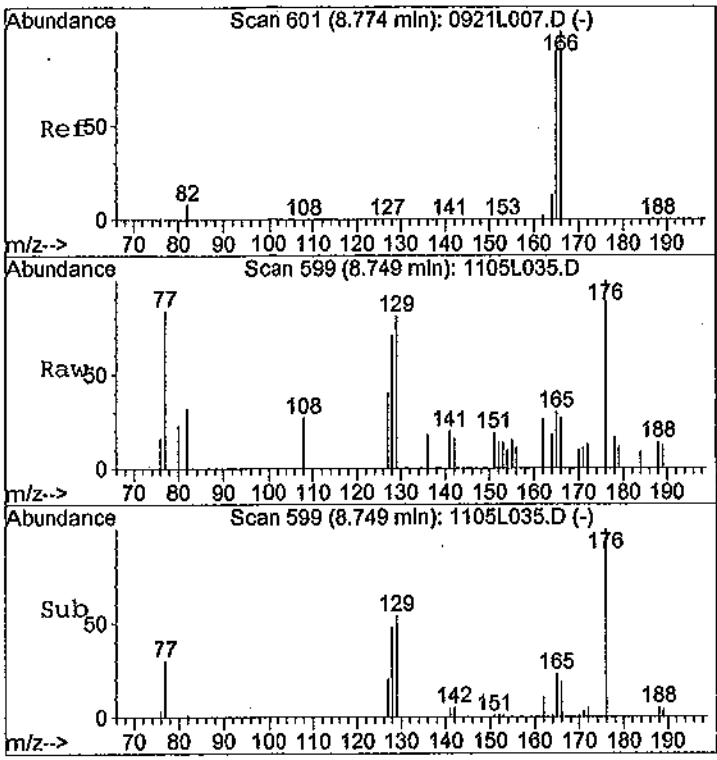


#9
Acenaphthene
Concen: 0.17136 ppb
RT: 8.15 min Scan# 549
Delta R.T. -0.00 min
Lab File: 1105L035.D
Acq: 5 Nov 11 23:11

Tgt Ion:154 Resp: 168

Ion	Ratio	Lower	Upper
154	100		
153	95.1	72.0	133.8
152	39.2	32.1	59.5

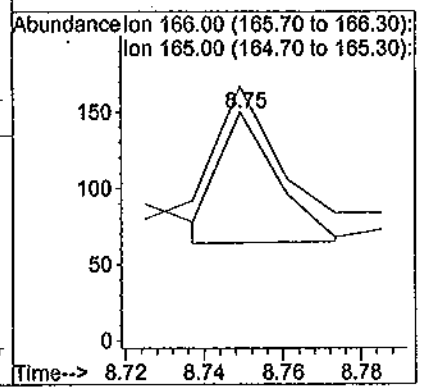




#10
Fluorene
Concen: 0.08203 ppb
RT: 8.75 min Scan# 599
Delta R.T. -0.01 min
Lab File: 1105L035.D
Acq: 5 Nov 11 23:11

Tgt Ion:166 Resp: 88

Ion	Ratio	Lower	Upper
166	100		
165	101.2	65.6	121.8



EPA 8270D SIM

Environet, Inc.
650 Iwilei Rd, #204
Honolulu, HI 96817

APPL Inc.
908 North Temperance Avenue
Clovis, CA 93611

Attn: Stacy Fineran
Project: RED HILL/1022-024
Sample ID: ES047
Sample Collection Date: 10/24/11

ARF: 66102
APPL ID: AY49334
QCG: #SIMHC-111031A-161019

Method	Analyte	Result	LOQ	LOD	DL	Units	Extraction Date	Analysis Date
8270D-SIM	1-METHYLNAPHTHALENE	0.62	0.2	0.12	0.06	ug/L	10/31/11	11/06/11
8270D-SIM	2-METHYLNAPHTHALENE	0.20	0.2	0.12	0.06	ug/L	10/31/11	11/06/11
8270D-SIM	ACENAPHTHENE	0.16 J	0.2	0.12	0.06	ug/L	10/31/11	11/06/11
8270D-SIM	ACENAPHTHYLENE	0.12 U	0.2	0.12	0.06	ug/L	10/31/11	11/06/11
8270D-SIM	ANTHRACENE	0.10 U	0.2	0.10	0.05	ug/L	10/31/11	11/06/11
8270D-SIM	BENZO(A)ANTHRACENE	0.14 U	0.2	0.14	0.07	ug/L	10/31/11	11/06/11
8270D-SIM	BENZO(A)PYRENE	0.14 U	0.2	0.14	0.07	ug/L	10/31/11	11/06/11
8270D-SIM	BENZO(B)FLUORANTHENE	0.12 U	0.2	0.12	0.06	ug/L	10/31/11	11/06/11
8270D-SIM	BENZO(GHI)PERYLENE	0.16 U	0.2	0.16	0.08	ug/L	10/31/11	11/06/11
8270D-SIM	BENZO(K)FLUORANTHENE	0.14 U	0.2	0.14	0.07	ug/L	10/31/11	11/06/11
8270D-SIM	CHRYSENE	0.10 U	0.2	0.10	0.05	ug/L	10/31/11	11/06/11
8270D-SIM	DIBENZ(A,H)ANTHRACENE	0.10 U	0.2	0.10	0.05	ug/L	10/31/11	11/06/11
8270D-SIM	FLUORANTHENE	0.16 U	0.2	0.16	0.08	ug/L	10/31/11	11/06/11
8270D-SIM	FLUORENE	0.083 J	0.2	0.12	0.06	ug/L	10/31/11	11/06/11
8270D-SIM	INDENO(1,2,3-CD)PYRENE	0.14 U	0.2	0.14	0.07	ug/L	10/31/11	11/06/11
8270D-SIM	NAPHTHALENE	1.0	0.2	0.10	0.05	ug/L	10/31/11	11/06/11
8270D-SIM	PHENANTHRENE	0.14 U	0.2	0.14	0.07	ug/L	10/31/11	11/06/11
8270D-SIM	PYRENE	0.16 U	0.2	0.16	0.08	ug/L	10/31/11	11/06/11
8270D-SIM	SURROGATE: 2-FLUORBIPHENYL (S)	53.4	50-110			%	10/31/11	11/06/11
8270D-SIM	SURROGATE: NITROBENZENE-D5 (S)	65.4	40-110			%	10/31/11	11/06/11
8270D-SIM	SURROGATE: TERPHENYL-D14 (S)	61.4	50-135			%	10/31/11	11/06/11

J = Estimated value.

Quant Method: SIM2.M
Run #: 1105L038
Instrument: Linus
Sequence: L111027
Dilution Factor: 1
Initials: LF

Printed: 11/09/11 3:59:14 PM
APPL-F1-SC-NoMC-REG MDLs

Quantitation Report (QT Reviewed)

Data File : M:\LINUS\DATA\L111027\1105L038.D Vial: 38
Acq On : 6 Nov 11 00:26 Operator: LF
Sample : AY49334W29 1/1050 Inst : Linus
Misc : Multiplr: 0.95

Quant Time: Nov 9 9:19 2011 Quant Results File: SIM2.RES

Quant Method : M:\LINUS\DATA\L111027\SIM2.M (RTE Integrator)
Title : EPA 8270C
Last Update : Wed Nov 02 15:56:51 2011
Response via : Initial Calibration
DataAcq Meth : 87SIMAQ

Internal Standards	R.T.	QIon	Response	Conc Units	Dev(Min)
1) Napthalene-D8(IS)	6.11	136	2328	2.50000 ppb	-0.01
6) Acenaphthene-D10(IS)	8.11	164	1134	2.50000 ppb	0.00
11) Phenanthrene-D10(IS)	9.84	188	1938	2.50000 ppb	-0.01
15) Chrysene-D12(IS)	12.94	240	2444	2.50000 ppb	0.01
21) Perylene-D12(IS)	14.57	264	2139	2.50000 ppb	0.01
System Monitoring Compounds					
2) Surrogate Recovery (NBZ)	5.41	82	545	1.24491 ppb	-0.04
Spiked Amount 1.905			Recovery = 65.363%		
7) Surrogate Recovery (FBP)	7.35	172	1080	1.01746 ppb	0.00
Spiked Amount 1.905			Recovery = 53.393%		
17) Surrogate Recovery (TPH)	11.70	244	1293	1.17013 ppb	-0.01
Spiked Amount 1.905			Recovery = 61.425%		
Target Compounds					Qvalue
3) Naphthalene	6.13	128	1754	1.02993 ppb #	54
4) 2-Methylnaphthalene	6.93	142	176	0.20154 ppb	94
5) 1-Methylnaphthalene	7.02	142	620	0.61508 ppb	93
9) Acenaphthene	8.15	154	143	0.15769 ppb	95
10) Fluorene	8.75	166	82	0.08264 ppb	92

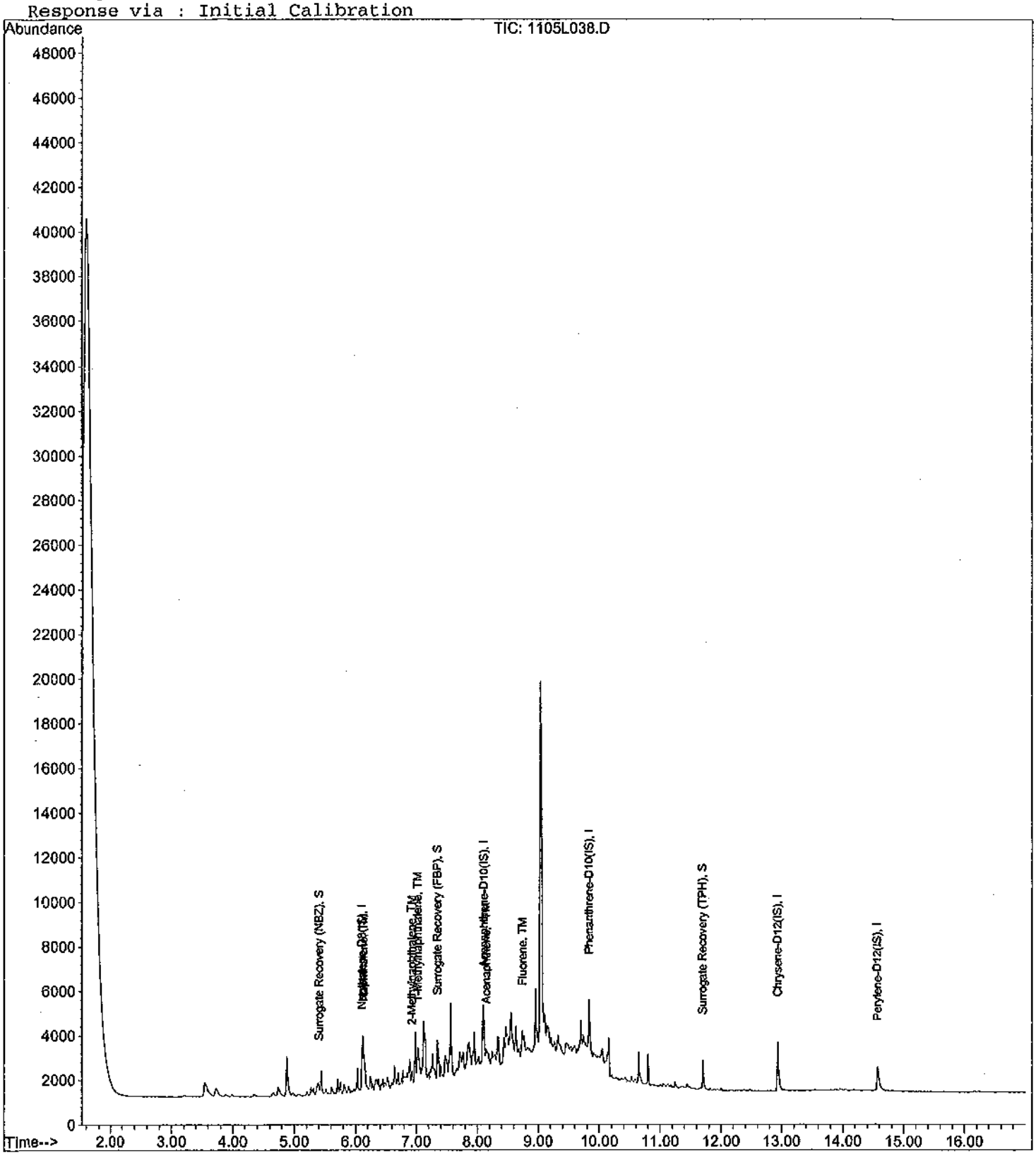
(#) = qualifier out of range (m) = manual integration

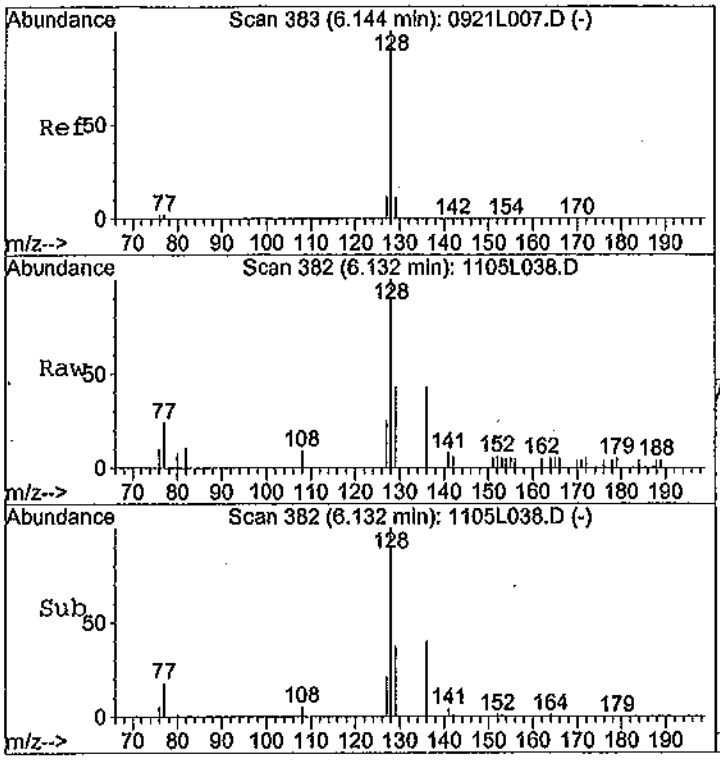
Quantitation Report

Data File : M:\LINUS\DATA\L111027\1105L038.D Vial: 38
Acq On : 6 Nov 11 00:26 Operator: LF
Sample : AY49334W29 1/1050 Inst : Linus
Misc : Multiplr: 0.95

Quant Time: Nov 9 9:19 2011 Quant Results File: SIM2.RES

Method : M:\LINUS\DATA\L111027\SIM2.M (RTE Integrator)
Title : EPA 8270C
Last Update : Tue Nov 08 16:22:04 2011
Response via : Initial Calibration

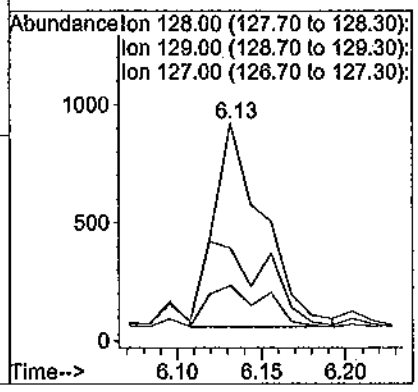


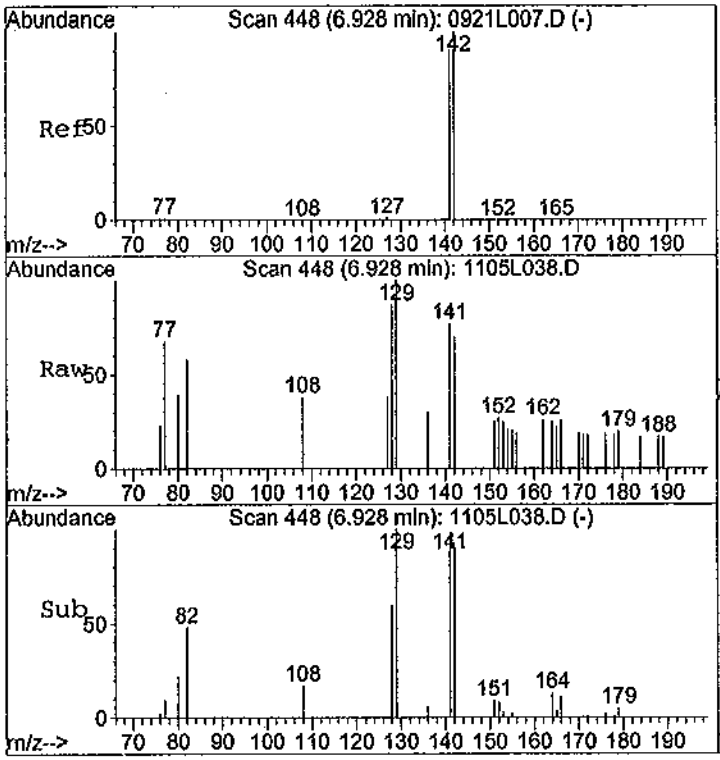


#3
Naphthalene
Concen: 1.02993 ppb
RT: 6.13 min Scan# 382
Delta R.T. -0.01 min
Lab File: 1105L038.D
Acq: 6 Nov 11 00:26

Tgt Ion:128 Resp: 1754

Ion	Ratio	Lower	Upper
128	100		
129	39.0	7.7	14.3#
127	20.8	8.6	16.0#

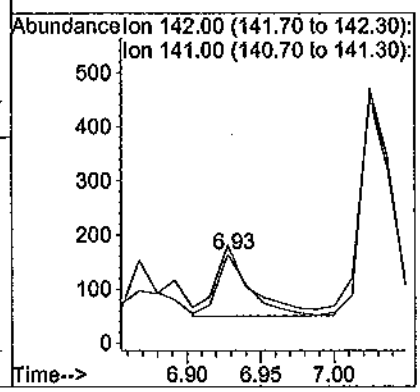


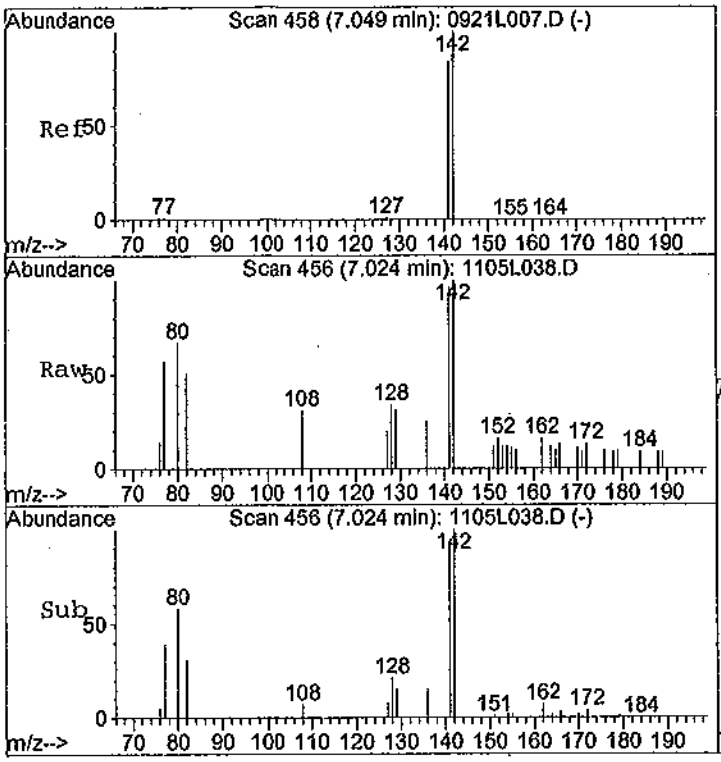


#4
2-Methylnaphthalene
Concen: 0.20154 ppb
RT: 6.93 min Scan# 448
Delta R.T. -0.04 min
Lab File: 1105L038.D
Acq: 6 Nov 11 00:26

Tgt Ion:142 Resp: 176

Ion	Ratio	Lower	Upper
142	100		
141	104.5	68.9	127.9

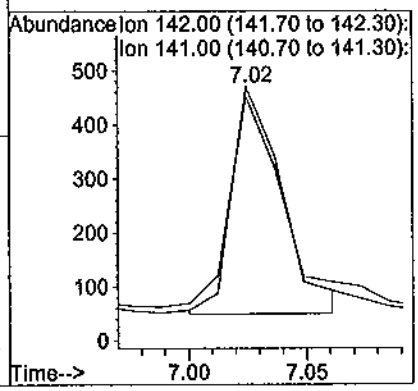


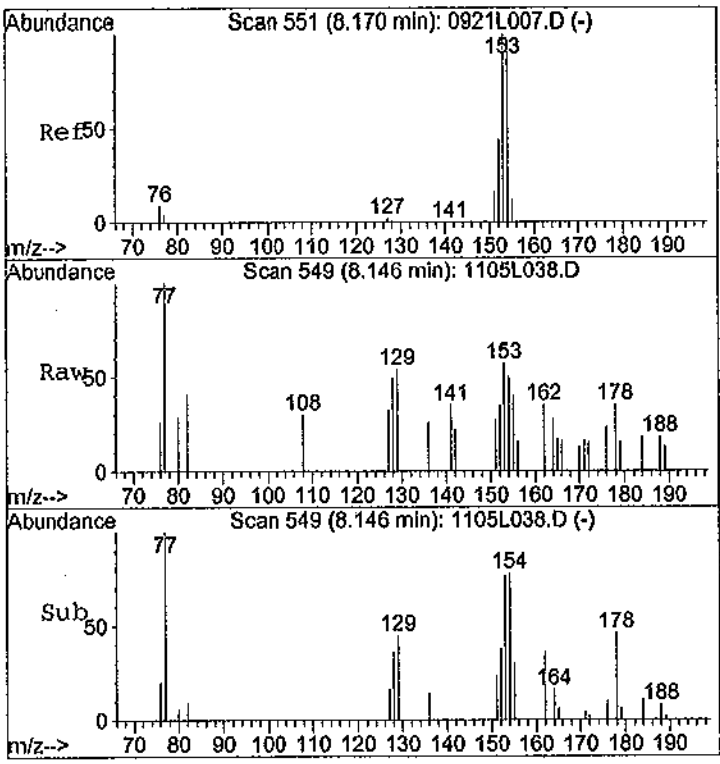


#5
1-Methylnaphthalene
Concen: 0.61508 ppb
RT: 7.02 min Scan# 456
Delta R.T. -0.02 min
Lab File: 1105L038.D
Acq: 6 Nov 11 00:26

Tgt Ion:142 Resp: 620

Ion	Ratio	Lower	Upper
142	100		
141	93.0	70.0	130.0

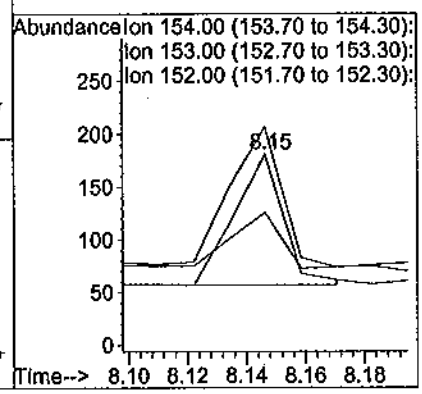


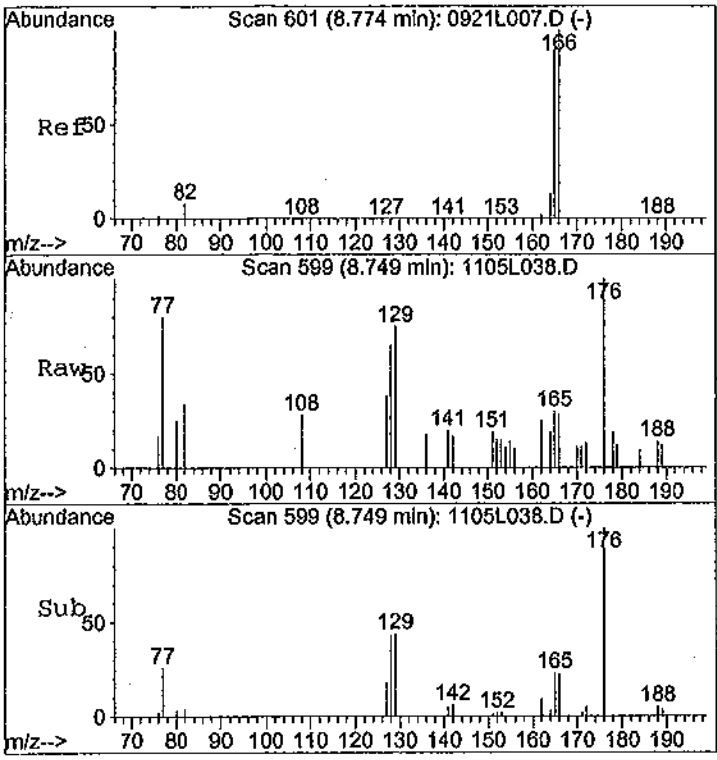


#9
Acenaphthene
Concen: 0.15769 ppb
RT: 8.15 min Scan# 549
Delta R.T. 0.00 min
Lab File: 1105L038.D
Acq: 6 Nov 11 00:26

Tgt Ion:154 Resp: 143

Ion	Ratio	Lower	Upper
154	100		
153	107.3	72.0	133.8
152	41.9	32.1	59.5

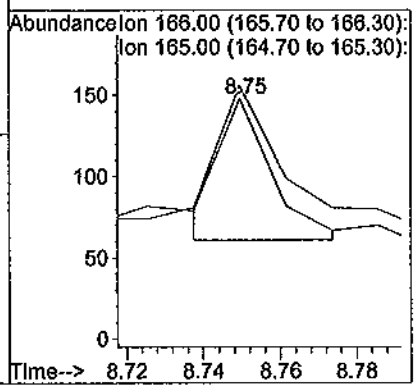




#10
Fluorene
Concen: 0.08264 ppb
RT: 8.75 min Scan# 599
Delta R.T. -0.01 min
Lab File: 1105L038.D
Acq: 6 Nov 11 00:26

Tgt Ion:166 Resp: 82

Ion	Ratio	Lower	Upper
166	100		
165	105.4	65.6	121.8



EPA 8270D SIM

Environet, Inc.
650 Iwilei Rd, #204
Honolulu, HI 96817

APPL Inc.
908 North Temperance Avenue
Clovis, CA 93611

Attn: Stacy Fineran
Project: RED HILL/1022-024
Sample ID: ES049
Sample Collection Date: 10/24/11

ARF: 66102
APPL ID: AY49336
QCG: #SIMHC-111031A-161019

Method	Analyte	Result	LOQ	LOD	DL	Units	Extraction Date	Analysis Date
8270D-SIM	1-METHYLNAPHTHALENE	0.12 U	0.2	0.12	0.06	ug/L	10/31/11	11/06/11
8270D-SIM	2-METHYLNAPHTHALENE	0.12 U	0.2	0.12	0.06	ug/L	10/31/11	11/06/11
8270D-SIM	ACENAPHTHENE	0.12 U	0.2	0.12	0.06	ug/L	10/31/11	11/06/11
8270D-SIM	ACENAPHTHYLENE	0.12 U	0.2	0.12	0.06	ug/L	10/31/11	11/06/11
8270D-SIM	ANTHRACENE	0.10 U	0.2	0.10	0.05	ug/L	10/31/11	11/06/11
8270D-SIM	BENZO(A)ANTHRACENE	0.14 U	0.2	0.14	0.07	ug/L	10/31/11	11/06/11
8270D-SIM	BENZO(A)PYRENE	0.14 U	0.2	0.14	0.07	ug/L	10/31/11	11/06/11
8270D-SIM	BENZO(B)FLUORANTHENE	0.12 U	0.2	0.12	0.06	ug/L	10/31/11	11/06/11
8270D-SIM	BENZO(GHI)PERYLENE	0.16 U	0.2	0.16	0.08	ug/L	10/31/11	11/06/11
8270D-SIM	BENZO(K)FLUORANTHENE	0.14 U	0.2	0.14	0.07	ug/L	10/31/11	11/06/11
8270D-SIM	CHRYSENE	0.10 U	0.2	0.10	0.05	ug/L	10/31/11	11/06/11
8270D-SIM	DIBENZ(A,H)ANTHRACENE	0.10 U	0.2	0.10	0.05	ug/L	10/31/11	11/06/11
8270D-SIM	FLUORANTHENE	0.16 U	0.2	0.16	0.08	ug/L	10/31/11	11/06/11
8270D-SIM	FLUORENE	0.12 U	0.2	0.12	0.06	ug/L	10/31/11	11/06/11
8270D-SIM	INDENO(1,2,3-CD)PYRENE	0.14 U	0.2	0.14	0.07	ug/L	10/31/11	11/06/11
8270D-SIM	NAPHTHALENE	0.10 U	0.2	0.10	0.05	ug/L	10/31/11	11/06/11
8270D-SIM	PHENANTHRENE	0.14 U	0.2	0.14	0.07	ug/L	10/31/11	11/06/11
8270D-SIM	PYRENE	0.16 U	0.2	0.16	0.08	ug/L	10/31/11	11/06/11
8270D-SIM	SURROGATE: 2-FLUORBIPHENYL (S)	65.0	50-110			%	10/31/11	11/06/11
8270D-SIM	SURROGATE: NITROBENZENE-D5 (S)	64.4	40-110			%	10/31/11	11/06/11
8270D-SIM	SURROGATE: TERPHENYL-D14 (S)	51.9	50-135			%	10/31/11	11/06/11

Quant Method: SIM2.M
Run #: 1105L039
Instrument: Linus
Sequence: L111027
Dilution Factor: 1
Initials: LF

Printed: 11/09/11 3:59:14 PM
APPL-F1-SC-NoMC-REG MDLs

Quantitation Report (QT Reviewed)

Data File : M:\LINUS\DATA\L111027\1105L039.D Vial: 39
Acq On : 6 Nov 11 00:51 Operator: LF
Sample : AY49336W10 1/1030 Inst : Linus
Misc : Multiplr: 0.97

Quant Time: Nov 9 9:20 2011 Quant Results File: SIM2.RES

Quant Method : M:\LINUS\DATA\L111027\SIM2.M (RTE Integrator)
Title : EPA 8270C
Last Update : Wed Nov 02 15:56:51 2011
Response via : Initial Calibration
DataAcq Meth : 87SIMAQ

Internal Standards	R.T.	QIon	Response	Conc Units	Dev(Min)
1) Napthalene-D8(IS)	6.12	136	2268	2.50000 ppb	0.00
6) Acenaphthene-D10(IS)	8.11	164	1052	2.50000 ppb	0.00
11) Phenanthrene-D10(IS)	9.86	188	1894	2.50000 ppb	0.01
15) Chrysene-D12(IS)	12.94	240	2526	2.50000 ppb	0.01
21) Perylene-D12(IS)	14.57	264	2131	2.50000 ppb	0.01
System Monitoring Compounds					
2) Surrogate Recovery (NBZ)	5.46	82	523	1.25007 ppb	0.01
Spiked Amount 1.942			Recovery = 64.375%		
7) Surrogate Recovery (FBP)	7.36	172	1220	1.26300 ppb	0.01
Spiked Amount 1.942			Recovery = 65.045%		
17) Surrogate Recovery (TPH)	11.71	244	1129	1.00774 ppb	0.00
Spiked Amount 1.942			Recovery = 51.912%		

Target Compounds Qvalue

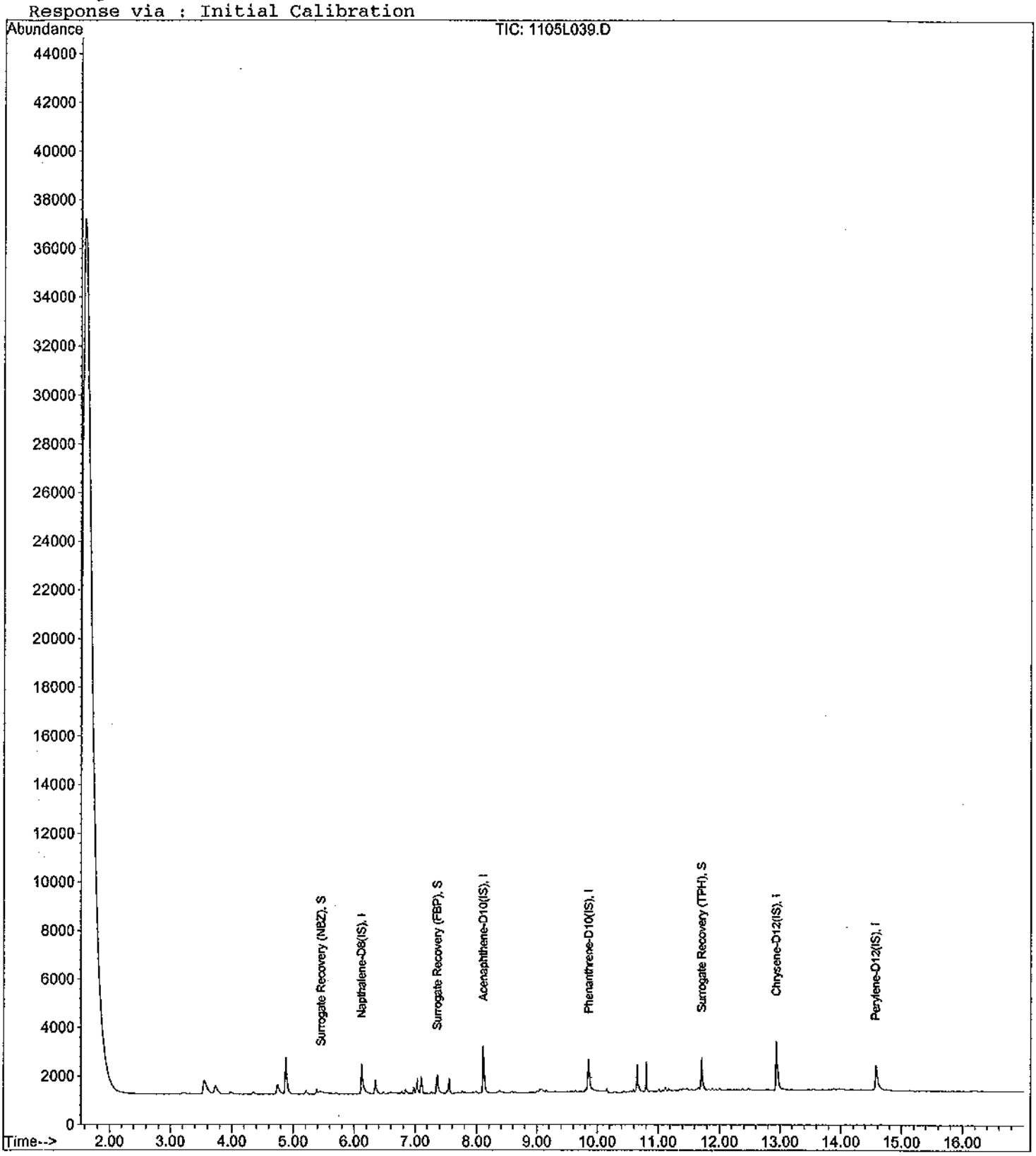
(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : M:\LINUS\DATA\L111027\1105L039.D Vial: 39
Acq On : 6 Nov 11 00:51 Operator: LF
Sample : AY49336W10 1/1030 Inst : Linus
Misc : Multiplr: 0.97

Quant Time: Nov 9 9:20 2011 Quant Results File: SIM2.RES

Method : M:\LINUS\DATA\L111027\SIM2.M (RTE Integrator)
Title : EPA 8270C
Last Update : Tue Nov 08 16:22:04 2011
Response via : Initial Calibration



EPA METHOD 8270
Polynuclear Aromatic Hydrocarbons
Calibration Data

APPL, INC.

EPA 8270C SIM

Form 6
Initial Calibration

Lab Name: APPL, Inc. SDG No: 6612

Case No: Initial Cal. Date: 10/27/11

Matrix: Instrument: Linus

Initials:

			1027L003.D	1027L004.D	1028L005.D	1028L006.D	1028L007.D	1028L008.D	1028L009.D	1028L010.D						
		Compound	0.1	0.2	0.5	1	5	10	50	100			Avg	%RSD		
1	I	Napthalene-D8(IS)														
2	S	Surrogate Recovery (NBZ)	0.4127	0.4673	0.4981	0.4588	0.3927	0.4106	0.4509	0.4908			0.45	8.7	S	
3	TM	Naphthalene	1.848	2.053	1.897	1.975	1.484	1.725	1.497	1.455			1.7	14	TM	
4	TM	2-Methylnaphthalene	0.8339	0.8429	0.8095	0.9618	0.8735	1.026	0.9052	0.8919			0.89	8.0	TM	
5	TM	1-Methylnaphthalene	1.006	1.140	1.127	1.262	0.9445	1.066	0.8694	0.8324			1.0	14	TM	
6	I	Acenaphthene-D10(IS)														
7	S	Surrogate Recovery (FBP)	2.266	2.373	2.477	2.606	2.184	2.112	1.948	1.864			2.2	11	S	
8	TM	Acenaphthylene	3.556	3.474	3.469	3.747	3.046	3.265	3.097	2.960			3.3	8.4	TM	
9	*TM	Acenaphthene	2.197	2.003	2.015	2.153	1.761	1.820	1.687	1.595			1.9	12	*TM	
10	TM	Fluorene	2.179	2.202	2.246	2.358	1.948	2.057	1.915	1.761			2.1	9.5	TM	
11	I	Phenanthrene-D10(IS)														
12	TM	Phenanthrene	1.851	1.670	1.765	1.775	1.470	1.719	1.377	1.248			1.6	13	TM	
13	TM	Anthracene	1.736	1.681	1.869	1.846	1.493	1.754	1.415	1.278			1.6	13	TM	
14	*TM	Fluoranthene	3.116	3.004	3.048	3.052	2.530	3.014	2.407	2.162			2.8	13	*TM	
15	I	Chrysene-D12(IS)														
16	TM	Pyrene	2.537	2.408	2.291	2.412	2.045	2.168	1.892	1.849			2.2	12	TM	
17	S	Surrogate Recovery (TPH)	1.147	1.155	1.070	1.235	1.097	1.067	0.9366	0.9044			1.1	10	S	
18	TM	Benz (a) anthracene	1.767	1.435	1.243	1.561	1.267	1.372	1.372	1.578			1.4	12	TM	
19	TM	Chrysene	2.135	2.066	2.037	2.114	1.981	2.084	1.655	1.438			1.9	13	TM	
20	TM	Indeno (1,2,3-cd) pyrene	1.825	1.519	1.284	1.532	1.378	1.521	1.460	1.500			1.5	10	TM	
21	I	Perylene-D12(IS)														
22	TM	Benzo (b) fluoranthene	1.711	1.652	1.826	1.857	1.789	1.866	1.765	1.625			1.8	5.2	TM	
23	TM	Benzo (k) fluoranthene	2.136	1.797	1.882	1.854	1.869	1.935	1.561	1.548			1.8	11	TM	
24	*TM	Benzo (a) pyrene	2.001	1.899	1.650	1.686	1.673	1.757	1.585	1.533			1.7	9.2	*TM	
25	TM	Dibenz (a,h) anthracene	1.773	1.577	1.175	1.438	1.356	1.477	1.399	1.379			1.4	12	TM	
26	TM	Benzo (g,h,i) perylene	1.410	1.754	1.503	1.639	1.449	1.558	1.444	1.446			1.5	7.8	TM	
27																
28																
29																
30																
31																
32																
33																
34																
35																

Quantitation Report (QT Reviewed)

Data File : M:\LINUS\DATA\L111027\1027L003.D Vial: 3
Acq On : 27 Oct 11 19:12 Operator: LF
Sample : 0.1ug/ml PAH 10-27-11 Inst : Linus
Misc : Multiplr: 1.00

Quant Time: Oct 30 11:15 2011 Quant Results File: SIM2.RES

Quant Method : M:\LINUS\DATA\L111028\SIM2.M (RTE Integrator)
Title : EPA 8270C
Last Update : Sun Oct 30 10:57:42 2011
Response via : Initial Calibration
DataAcq Meth : 87SIMAQ

Internal Standards	R.T.	QIon	Response	Conc Units		Dev(Min)
1) Napthalene-D8(IS)	6.14	136	2908	2.50000 ppb		0.02
6) Acenaphthene-D10(IS)	8.12	164	1434	2.50000 ppb		0.01
11) Phenanthrene-D10(IS)	9.87	188	2391	2.50000 ppb		0.02
15) Chrysene-D12(IS)	12.95	240	2986	2.50000 ppb		0.02
21) Perylene-D12(IS)	14.58	264	2411	2.50000 ppb		0.02
System Monitoring Compounds						
2) Surrogate Recovery (NBZ)	5.61	82	48	0.74306 ppb		0.19
Spiked Amount 2.000			Recovery	= 37.150%		
7) Surrogate Recovery (FBP)	7.40	172	130	0.09815 ppb		0.05
Spiked Amount 2.000			Recovery	= 4.900%		
17) Surrogate Recovery (TPH)	11.74	244	137	0.09107 ppb		0.02
Spiked Amount 2.000			Recovery	= 4.550%		
Target Compounds						Qvalue
3) Naphthalene	6.17	128	215	0.10425 ppb		93
4) 2-Methylnaphthalene	7.01	142	97	0.09198 ppb		99
5) 1-Methylnaphthalene	7.08	142	117	0.09071 ppb		97
8) Acenaphthylene	7.99	152	204	0.10524 ppb		99
9) Acenaphthene	8.16	154	126	0.11351 ppb		94
10) Fluorene	8.81	166	125	0.10297 ppb		98
12) Phenanthrene	9.90	178	177	0.11216 ppb		95
13) Anthracene	9.99	178	166	0.10145 ppb		95
14) Fluoranthene	11.30	202	298	0.10883 ppb	#	90
16) Pyrene	11.56	202	303	0.11040 ppb		99
18) Benz (a) anthracene	12.95	228	211	0.11702 ppb		96
19) Chrysene	12.98	228	255	0.09385 ppb		98
20) Indeno (1,2,3-cd) pyrene	16.19	276	218	0.11665 ppb	#	93
22) Benzo (b) fluoranthene	14.15	252	165	0.09422 ppb	#	95
23) Benzo (k) fluoranthene	14.19	252	206	0.11693 ppb		65
24) Benzo (a) pyrene	14.54	252	193	0.11081 ppb		95
25) Dibenz (a,h) anthracene	16.17	278	171	0.11827 ppb		92
26) Benzo (g,h,i) perylene	16.64	276	136	0.08955 ppb	#	89

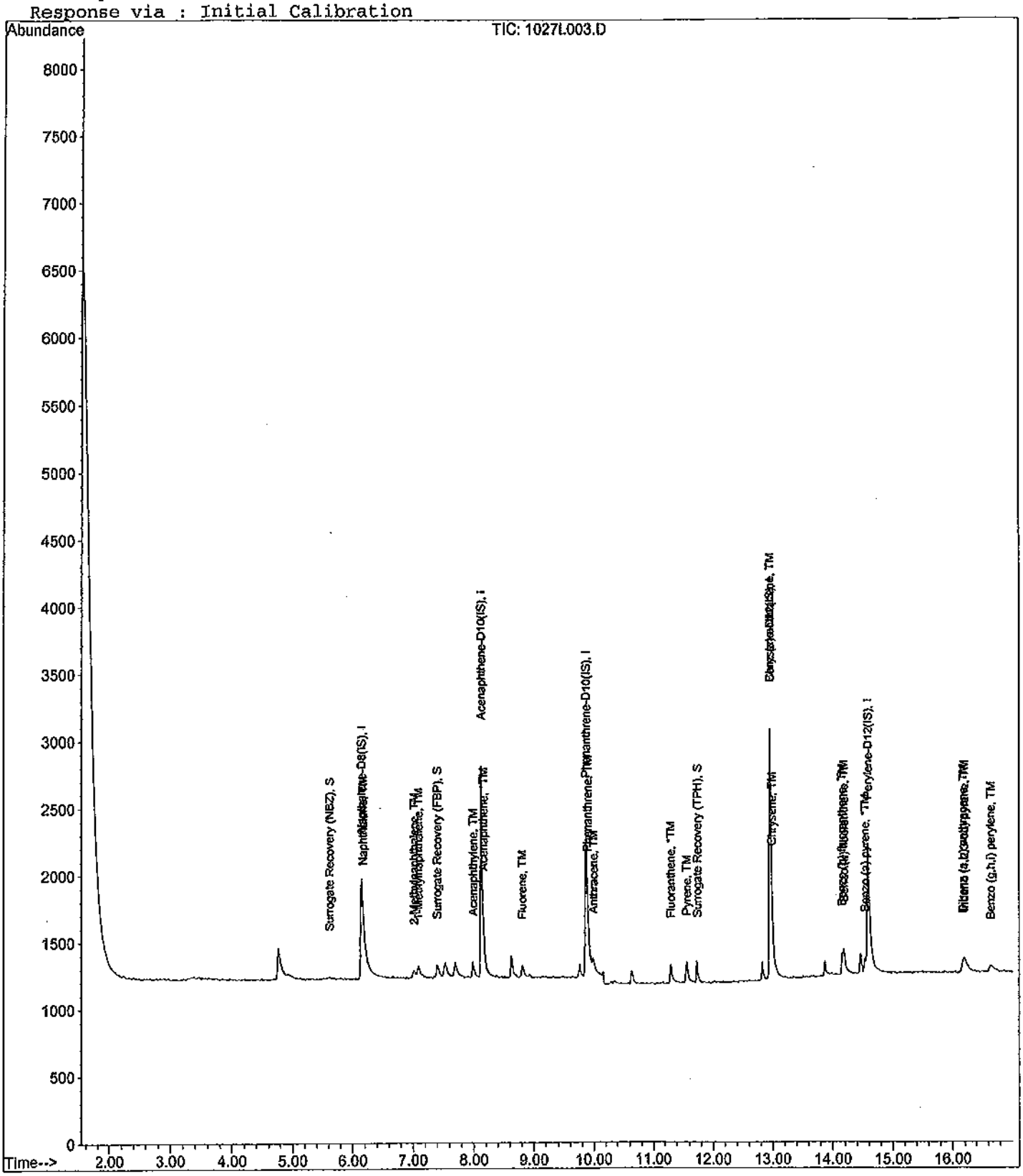
(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : M:\LINUS\DATA\L111027\1027L003.D Vial: 3
Acq On : 27 Oct 11 19:12 Operator: LF
Sample : 0.1ug/ml PAH 10-27-11 Inst : Linus
Misc : Multiplr: 1.00

Quant Time: Oct 30 11:15 2011 Quant Results File: SIM2.RES

Method : M:\LINUS\DATA\L111027\SIM2.M (RTE Integrator)
Title : EPA 8270C
Last Update : Tue Nov 01 17:14:29 2011
Response via : Initial Calibration



Quantitation Report (QT Reviewed)

Data File : M:\LINUS\DATA\L111027\1027L004.D
Acq On : 27 Oct 11 19:38
Sample : 0.2ug/ml PAH
Misc :

Vial: 4
Operator: LF
Inst : Linus
Multiplr: 1.00

Quant Time: Oct 30 11:13 2011 Quant Results File: SIM2.RES

Quant Method : M:\LINUS\DATA\L111028\SIM2.M (RTE Integrator)
Title : EPA 8270C
Last Update : Sun Oct 30 10:57:42 2011
Response via : Initial Calibration
DataAcq Meth : 87SIMAQ

Internal Standards	R.T.	QIon	Response	Conc Units		Dev(Min)
1) Napthalene-D8(IS)	6.14	136	2862	2.50000 ppb		0.02
6) Acenaphthene-D10(IS)	8.12	164	1317	2.50000 ppb		0.01
11) Phenanthrene-D10(IS)	9.87	188	2305	2.50000 ppb		0.02
15) Chrysene-D12(IS)	12.95	240	2814	2.50000 ppb		0.02
21) Perylene-D12(IS)	14.58	264	2323	2.50000 ppb		0.02
System Monitoring Compounds						
2) Surrogate Recovery (NBZ)	5.60	82	107	0.84083 ppb		0.18
Spiked Amount 2.000			Recovery = 42.050%			
7) Surrogate Recovery (FBP)	7.40	172	250	0.20995 ppb		0.05
Spiked Amount 2.000			Recovery = 10.500%			
17) Surrogate Recovery (TPH)	11.72	244	260	0.18421 ppb		0.01
Spiked Amount 2.000			Recovery = 9.200%			
Target Compounds						Qvalue
3) Naphthalene	6.17	128	470	0.23025 ppb		94
4) 2-Methylnaphthalene	7.00	142	193	0.18513 ppb		92
5) 1-Methylnaphthalene	7.07	142	261	0.20451 ppb		98
8) Acenaphthylene	7.99	152	366	0.20677 ppb		98
9) Acenaphthene	8.16	154	211	0.20826 ppb		87
10) Fluorene	8.81	166	232	0.20927 ppb		99
12) Phenanthrene	9.90	178	308	0.20239 ppb		96
13) Anthracene	9.99	178	310	0.19992 ppb		95
14) Fluoranthene	11.29	202	554	0.20981 ppb		95
16) Pyrene	11.55	202	542	0.21034 ppb	#	91
18) Benz (a) anthracene	12.95	228	323	0.19084 ppb		97
19) Chrysene	12.98	228	465	0.18296 ppb		99
20) Indeno (1,2,3-cd) pyrene	16.17	276	342	0.19494 ppb	#	96
22) Benzo (b) fluoranthene	14.15	252	307	0.18266 ppb		97
23) Benzo (k) fluoranthene	14.19	252	334	0.18857 ppb		64
24) Benzo (a) pyrene	14.54	252	353	0.21468 ppb		96
25) Dibenz (a,h) anthracene	16.16	278	293	0.21252 ppb		92
26) Benzo (g,h,i) perylene	16.64	276	326	0.22362 ppb		88

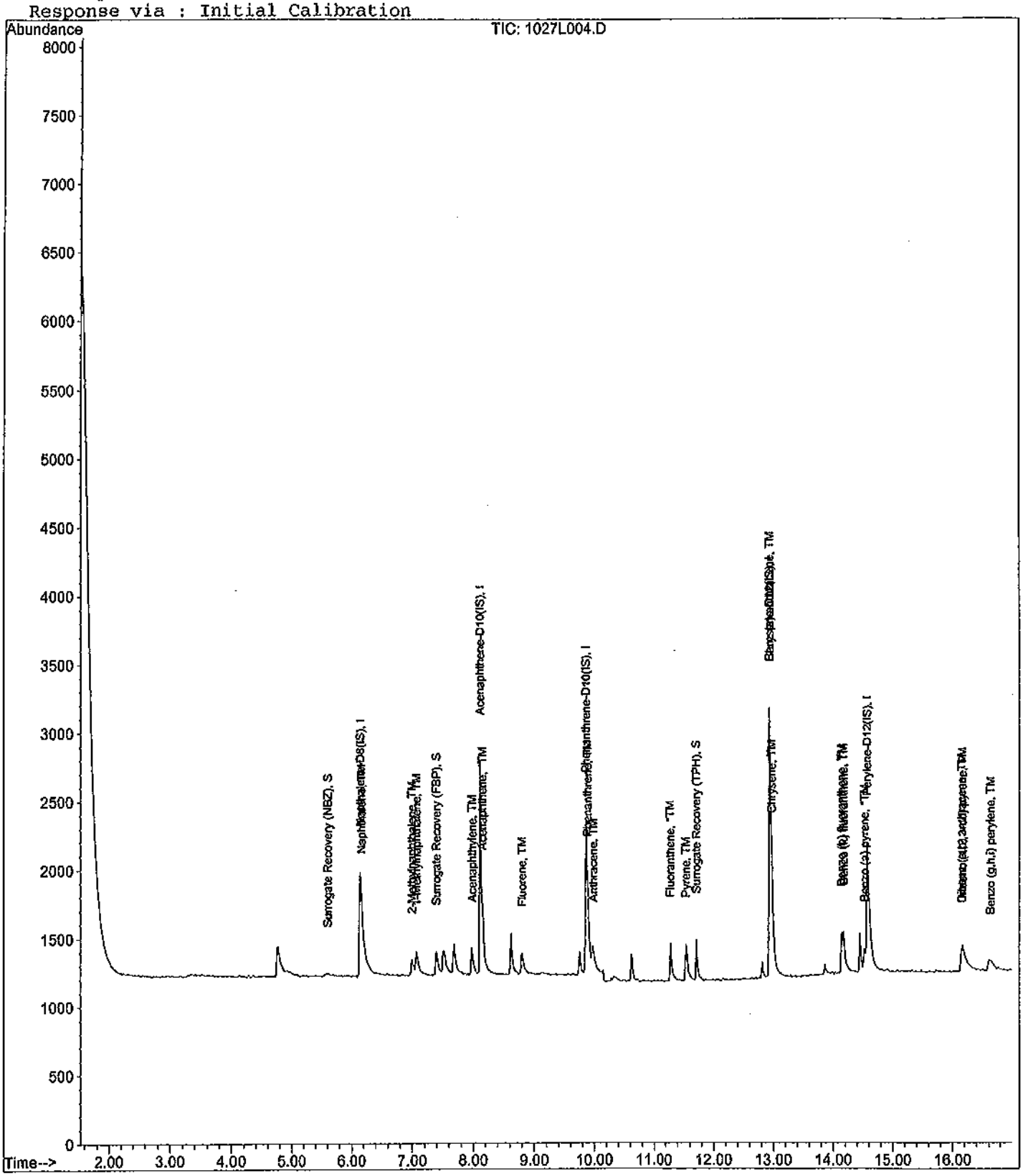
(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : M:\LINUS\DATA\L111027\1027L004.D Vial: 4
Acq On : 27 Oct 11 19:38 Operator: LF
Sample : 0.2ug/ml PAH Inst : Linus
Misc : Multiplr: 1.00

Quant Time: Oct 30 11:13 2011 Quant Results File: SIM2.RES

Method : M:\LINUS\DATA\L111027\SIM2.M (RTE Integrator)
Title : EPA 8270C
Last Update : Tue Nov 01 17:14:29 2011
Response via : Initial Calibration



Quantitation Report (QT Reviewed)

Data File : M:\LINUS\DATA\L111027\1028L005.D Vial: 5
Acq On : 28 Oct 11 11:07 Operator: LF
Sample : 0.5ug/ml PAH Inst : Linus
Misc : Multiplr: 1.00

Quant Time: Oct 30 11:12 2011 Quant Results File: SIM2.RES

Quant Method : M:\LINUS\DATA\L111028\SIM2.M (RTE Integrator)
Title : EPA 8270C
Last Update : Thu Sep 29 11:47:40 2011
Response via : Initial Calibration
DataAcq Meth : 87SIMAQ

Internal Standards	R.T.	QIon	Response	Conc Units		Dev(Min)
1) Napthalene-D8(IS)	6.14	136	2409	2.50000 ppb		0.02
6) Acenaphthene-D10(IS)	8.12	164	1104	2.50000 ppb		-0.01
11) Phenanthrene-D10(IS)	9.87	188	1819	2.50000 ppb		0.00
15) Chrysene-D12(IS)	12.95	240	2477	2.50000 ppb		-0.01
21) Perylene-D12(IS)	14.57	264	2043	2.50000 ppb		-0.02
System Monitoring Compounds						
2) Surrogate Recovery (NBZ)	5.60	82	240	1.15802 ppb		0.25
Spiked Amount 2.000			Recovery	= 57.900%		
7) Surrogate Recovery (FBP)	7.39	172	547	0.79241 ppb		0.01
Spiked Amount 2.000			Recovery	= 39.600%		
17) Surrogate Recovery (TPH)	11.74	244	530	0.66674 ppb		-0.02
Spiked Amount 2.000			Recovery	= 33.350%		
Target Compounds						Qvalue
3) Naphthalene	6.17	128	914	0.46769 ppb		98
4) 2-Methylnaphthalene	6.99	142	390	0.33945 ppb		96
5) 1-Methylnaphthalene	7.06	142	543	0.44086 ppb		95
8) Acenaphthylene	7.98	152	766	0.43771 ppb		99
9) Acenaphthene	8.16	154	445	0.43164 ppb		89
10) Fluorene	8.80	166	496	0.42124 ppb		99
12) Phenanthrene	9.90	178	642	0.38630 ppb		97
13) Anthracene	9.98	178	680	0.37229 ppb		95
14) Fluoranthene	11.29	202	1109	0.36672 ppb		96
16) Pyrene	11.55	202	1135	0.35574 ppb		97
18) Benz (a) anthracene	12.95	228	616	0.34309 ppb		98
19) Chrysene	12.98	228	1009	0.43128 ppb		99
20) Indeno (1,2,3-cd) pyrene	16.15	276	636	0.45186 ppb	#	96
22) Benzo (b) fluoranthene	14.14	252	746	0.48527 ppb		98
23) Benzo (k) fluoranthene	14.17	252	769	0.37285 ppb		98
24) Benzo (a) pyrene	14.52	252	674	0.41516 ppb		94
25) Dibenz (a,h) anthracene	16.14	278	480	0.46345 ppb		95
26) Benzo (g,h,i) perylene	16.59	276	614	0.46797 ppb		92

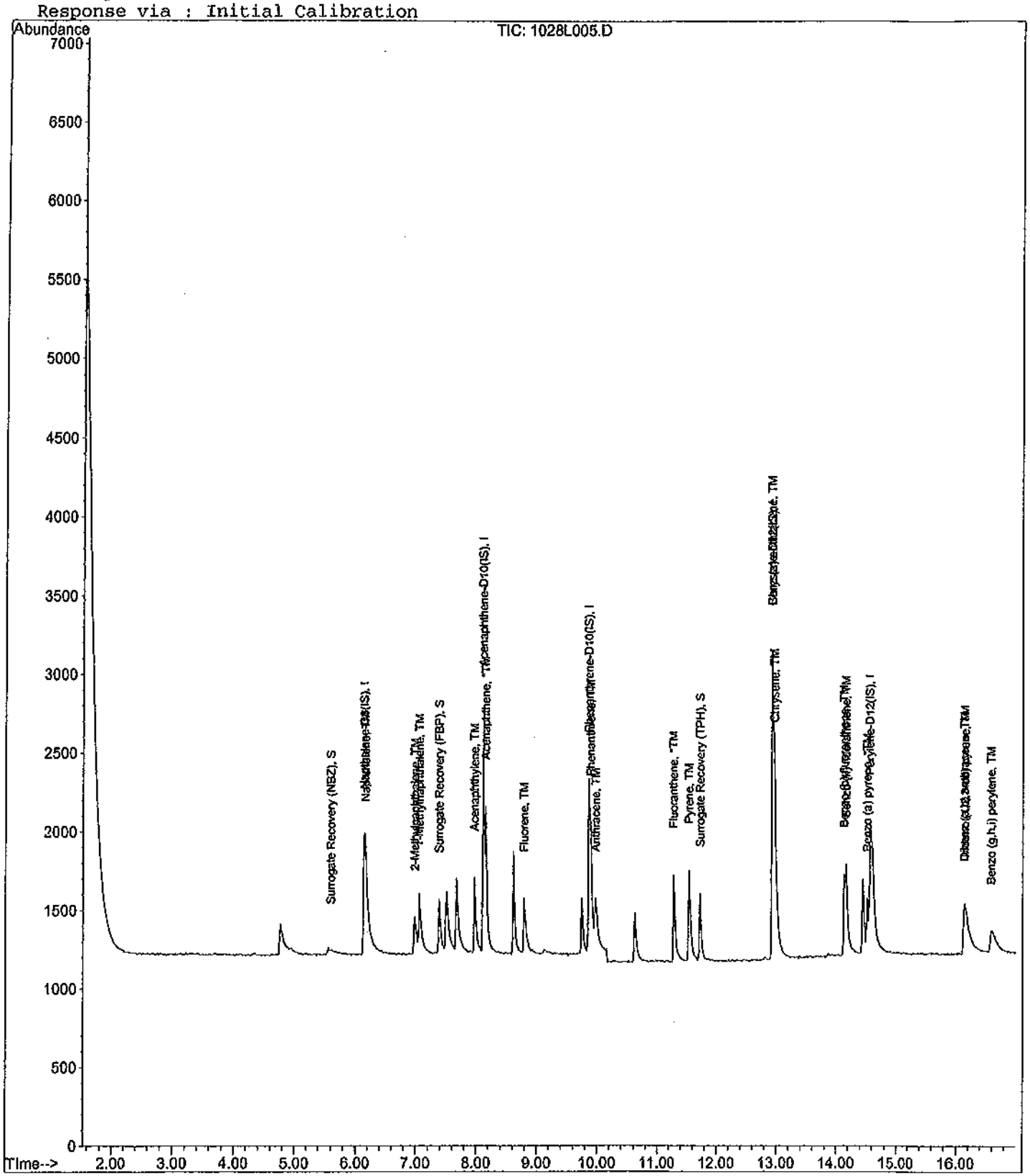
(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : M:\LINUS\DATA\L111027\1028L005.D Vial: 5
Acq On : 28 Oct 11 11:07 Operator: LF
Sample : 0.5ug/ml PAH Inst : Linus
Misc : Multiplr: 1.00

Quant Time: Oct 30 11:12 2011 Quant Results File: SIM2.RES

Method : M:\LINUS\DATA\L111027\SIM2.M (RTE Integrator)
Title : EPA 8270C
Last Update : Tue Nov 01 17:14:29 2011
Response via : Initial Calibration



Quantitation Report (QT Reviewed)

Data File : M:\LINUS\DATA\L111027\1028L006.D Vial: 6
Acq On : 28 Oct 11 11:32 Operator: LF
Sample : 1.0ug/ml PAH Inst : Linus
Misc : Multiplr: 1.00

Quant Time: Oct 30 11:10 2011 Quant Results File: SIM2.RES

Quant Method : M:\LINUS\DATA\L111028\SIM2.M (RTE Integrator)
Title : EPA 8270C
Last Update : Sun Oct 30 10:38:04 2011
Response via : Initial Calibration
DataAcq Meth : 87SIMAQ

Internal Standards	R.T.	QIon	Response	Conc Units		Dev(Min)
1) Napthalene-D8(IS)	6.13	136	2381	2.50000 ppb		0.01
6) Acenaphthene-D10(IS)	8.12	164	1089	2.50000 ppb		-0.01
11) Phenanthrene-D10(IS)	9.86	188	1865	2.50000 ppb		-0.01
15) Chrysene-D12(IS)	12.95	240	2449	2.50000 ppb		-0.01
21) Perylene-D12(IS)	14.57	264	2032	2.50000 ppb		-0.02
System Monitoring Compounds						
2) Surrogate Recovery (NBZ)	5.54	82	437	1.90266 ppb		0.00
Spiked Amount 2.000			Recovery = 95.150%			
7) Surrogate Recovery (FBP)	7.37	172	1135	1.66686 ppb		0.00
Spiked Amount 2.000			Recovery = 83.350%			
17) Surrogate Recovery (TPH)	11.72	244	1210	1.53959 ppb		-0.04
Spiked Amount 2.000			Recovery = 77.000%			
Target Compounds						Qvalue
3) Naphthalene	6.16	128	1881	0.97382 ppb		98
4) 2-Methylnaphthalene	6.96	142	916	0.80665 ppb		94
5) 1-Methylnaphthalene	7.05	142	1202	0.98738 ppb		89
8) Acenaphthylene	7.96	152	1632	0.94540 ppb		98
9) Acenaphthene	8.16	154	938	0.92237 ppb		91
10) Fluorene	8.79	166	1027	0.88422 ppb		98
12) Phenanthrene	9.90	178	1324	0.77703 ppb		99
13) Anthracene	9.97	178	1377	0.73529 ppb		98
14) Fluoranthene	11.28	202	2277	0.73437 ppb	#	94
16) Pyrene	11.54	202	2363	0.74909 ppb		97
18) Benz (a) anthracene	12.94	228	1529	0.86133 ppb		99
19) Chrysene	12.97	228	2071	0.89534 ppb		99
20) Indeno (1,2,3-cd) pyrene	16.12	276	1501	1.07861 ppb	#	92
22) Benzo (b) fluoranthene	14.13	252	1509	0.98690 ppb	#	96
23) Benzo (k) fluoranthene	14.16	252	1507	0.73463 ppb		96
24) Benzo (a) pyrene	14.51	252	1370	0.84844 ppb		98
25) Dibenz (a,h) anthracene	16.12	278	1169	1.13481 ppb		97
26) Benzo (g,h,i) perylene	16.58	276	1332	1.02070 ppb		98

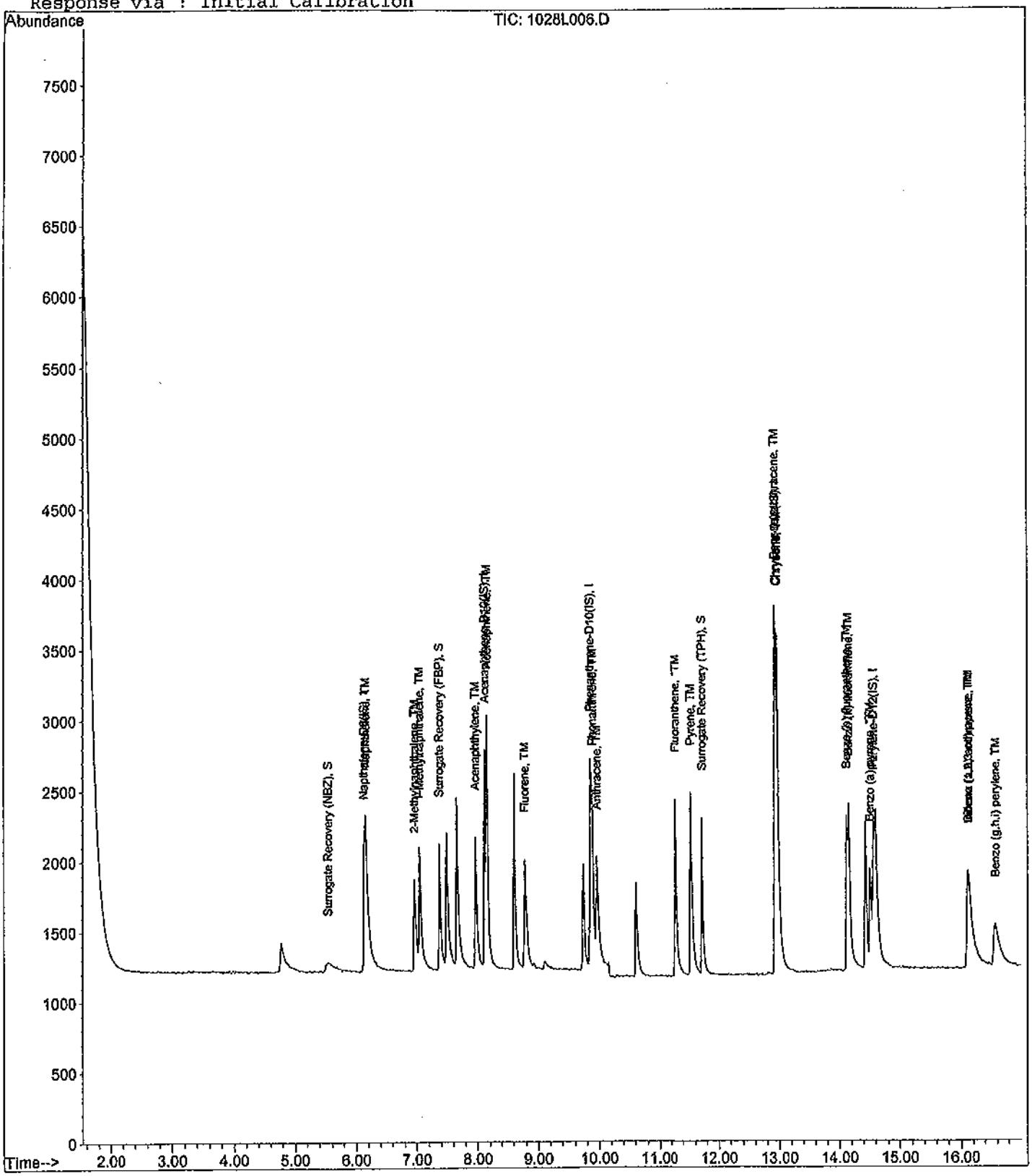
(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : M:\LINUS\DATA\L111027\1028L006.D Vial: 6
Acq On : 28 Oct 11 11:32 Operator: LF
Sample : 1.0ug/ml PAH Inst : Linus
Misc : Multiplr: 1.00

Quant Time: Oct 30 11:10 2011 Quant Results File: SIM2.RES

Method : M:\LINUS\DATA\L111027\SIM2.M (RTE Integrator)
Title : EPA 8270C
Last Update : Tue Nov 01 17:14:29 2011
Response via : Initial Calibration



Quantitation Report (QT Reviewed)

Data File : M:\LINUS\DATA\L111027\1028L007.D Vial: 7
Acq On : 28 Oct 11 11:58 Operator: LF
Sample : 5.0ug/ml PAH Inst : Linus
Misc : Multiplr: 1.00

Quant Time: Oct 30 10:40 2011 Quant Results File: SIM2.RES

Quant Method : M:\LINUS\DATA\L111028\SIM2.M (RTE Integrator)
Title : EPA 8270C
Last Update : Sun Oct 30 10:38:04 2011
Response via : Initial Calibration
DataAcq Meth : 87SIMAQ

Internal Standards	R.T.	QIon	Response	Conc Units		Dev(Min)
1) Napthalene-D8(IS)	6.12	136	2479	2.50000 ppb		0.00
6) Acenaphthene-D10(IS)	8.11	164	1083	2.50000 ppb		-0.02
11) Phenanthrene-D10(IS)	9.85	188	1851	2.50000 ppb		-0.02
15) Chrysene-D12(IS)	12.93	240	2378	2.50000 ppb		-0.04
21) Perylene-D12(IS)	14.56	264	1871	2.50000 ppb		-0.04
System Monitoring Compounds						
2) Surrogate Recovery (NBZ)	5.42	82	1947	7.24379 ppb		-0.12
Spiked Amount 2.000			Recovery	= 362.200%		
7) Surrogate Recovery (FBP)	7.35	172	4731	6.98644 ppb		-0.02
Spiked Amount 2.000			Recovery	= 349.300%		
17) Surrogate Recovery (TPH)	11.71	244	5216	6.83493 ppb		-0.05
Spiked Amount 2.000			Recovery	= 341.750%		
Target Compounds						Qvalue
3) Naphthalene	6.14	128	7358	3.65875 ppb		99
4) 2-Methylnaphthalene	6.93	142	4331	3.66320 ppb		98
5) 1-Methylnaphthalene	7.04	142	4683	3.69477 ppb		97
8) Acenaphthylene	7.95	152	6597	3.84274 ppb		100
9) Acenaphthene	8.15	154	3814	3.77124 ppb		92
10) Fluorene	8.76	166	4219	3.65257 ppb		99
12) Phenanthrene	9.87	178	5443	3.21854 ppb		98
13) Anthracene	9.94	178	5527	2.97363 ppb		99
14) Fluoranthene	11.26	202	9367	3.04387 ppb		98
16) Pyrene	11.51	202	9724	3.17462 ppb		97
18) Benz (a) anthracene	12.91	228	6027	3.49657 ppb		98
19) Chrysene	12.96	228	9422	4.19498 ppb		99
20) Indeno (1,2,3-cd) pyrene	16.06	276	6554	4.85029 ppb		95
22) Benzo (b) fluoranthene	14.10	252	6693	4.75397 ppb	#	96
23) Benzo (k) fluoranthene	14.14	252	6995	3.70332 ppb		99
24) Benzo (a) pyrene	14.49	252	6259	4.20974 ppb		98
25) Dibenz (a,h) anthracene	16.08	278	5075	5.35048 ppb		97
26) Benzo (g,h,i) perylene	16.51	276	5423	4.51321 ppb		98

(#) = qualifier out of range (m) = manual integration

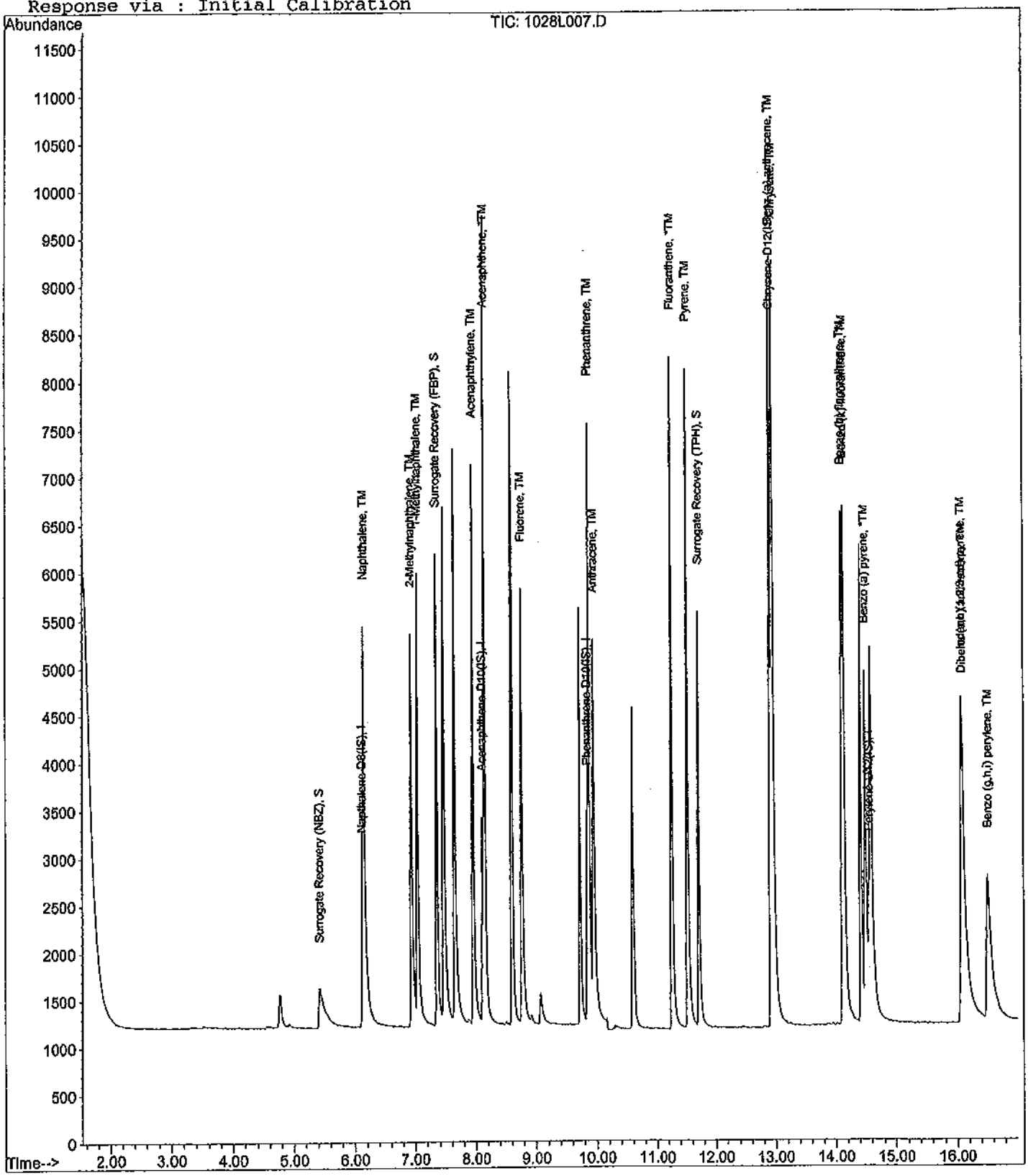
Quantitation Report

Data File : M:\LINUS\DATA\L111027\1028L007.D
Acq On : 28 Oct 11 11:58
Sample : 5.0ug/ml PAH
Misc :

Vial: 7
Operator: LF
Inst : Linus
Multiplr: 1.00

Quant Time: Oct 30 10:40 2011 Quant Results File: SIM2.RES

Method : M:\LINUS\DATA\L111027\SIM2.M (RTE Integrator)
Title : EPA 8270C
Last Update : Tue Nov 01 17:14:29 2011
Response via : Initial Calibration



Quantitation Report (QT Reviewed)

Data File : M:\LINUS\DATA\L111027\1028L008.D Vial: 8
Acq On : 28 Oct 11 12:23 Operator: LF
Sample : 10ug/ml PAH Inst : Linus
Misc : Multiplr: 1.00

Quant Time: Oct 30 10:41 2011 Quant Results File: SIM2.RES

Quant Method : M:\LINUS\DATA\L111028\SIM2.M (RTE Integrator)
Title : EPA 8270C
Last Update : Sun Oct 30 10:38:04 2011
Response via : Initial Calibration
DataAcq Meth : 87SIMAQ

Internal Standards	R.T.	QIon	Response	Conc Units		Dev(Min)
1) Napthalene-D8(IS)	6.12	136	2419	2.50000 ppb		0.00
6) Acenaphthene-D10(IS)	8.11	164	1154	2.50000 ppb		-0.02
11) Phenanthrene-D10(IS)	9.85	188	1800	2.50000 ppb		-0.02
15) Chrysene-D12(IS)	12.91	240	2580	2.50000 ppb		-0.05
21) Perylene-D12(IS)	14.55	264	2113	2.50000 ppb		-0.05
System Monitoring Compounds						
2) Surrogate Recovery (NBZ)	5.38	82	3973	14.84926 ppb		-0.16
Spiked Amount 2.000			Recovery = 742.450%			
7) Surrogate Recovery (FBP)	7.35	172	9747	13.50818 ppb		-0.02
Spiked Amount 2.000			Recovery = 675.400%			
17) Surrogate Recovery (TPH)	11.70	244	11014	13.30251 ppb		-0.06
Spiked Amount 2.000			Recovery = 665.150%			
Target Compounds						Qvalue
3) Naphthalene	6.13	128	16688	8.50390 ppb		99
4) 2-Methylnaphthalene	6.92	142	9930	8.60721 ppb		100
5) 1-Methylnaphthalene	7.02	142	10317	8.34175 ppb		92
8) Acenaphthylene	7.95	152	15071	8.23870 ppb		99
9) Acenaphthene	8.15	154	8403	7.79759 ppb		97
10) Fluorene	8.75	166	9496	7.71528 ppb		98
12) Phenanthrene	9.87	178	12375	7.52487 ppb		99
13) Anthracene	9.93	178	12631	6.98825 ppb		99
14) Fluoranthene	11.25	202	21698	7.25069 ppb	#	93
16) Pyrene	11.50	202	22373	6.73230 ppb	#	85
18) Benz (a) anthracene	12.91	228	14154	7.56854 ppb		100
19) Chrysene	12.95	228	21503	8.82425 ppb		99
20) Indeno (1,2,3-cd) pyrene	16.03	276	15698	10.70773 ppb	#	96
22) Benzo (b) fluoranthene	14.09	252	15772	9.91966 ppb		96
23) Benzo (k) fluoranthene	14.13	252	16351	7.66517 ppb		98
24) Benzo (a) pyrene	14.48	252	14853	8.84584 ppb		98
25) Dibenz (a,h) anthracene	16.05	278	12481	11.65147 ppb		96
26) Benzo (g,h,i) perylene	16.47	276	13167	9.70302 ppb		97

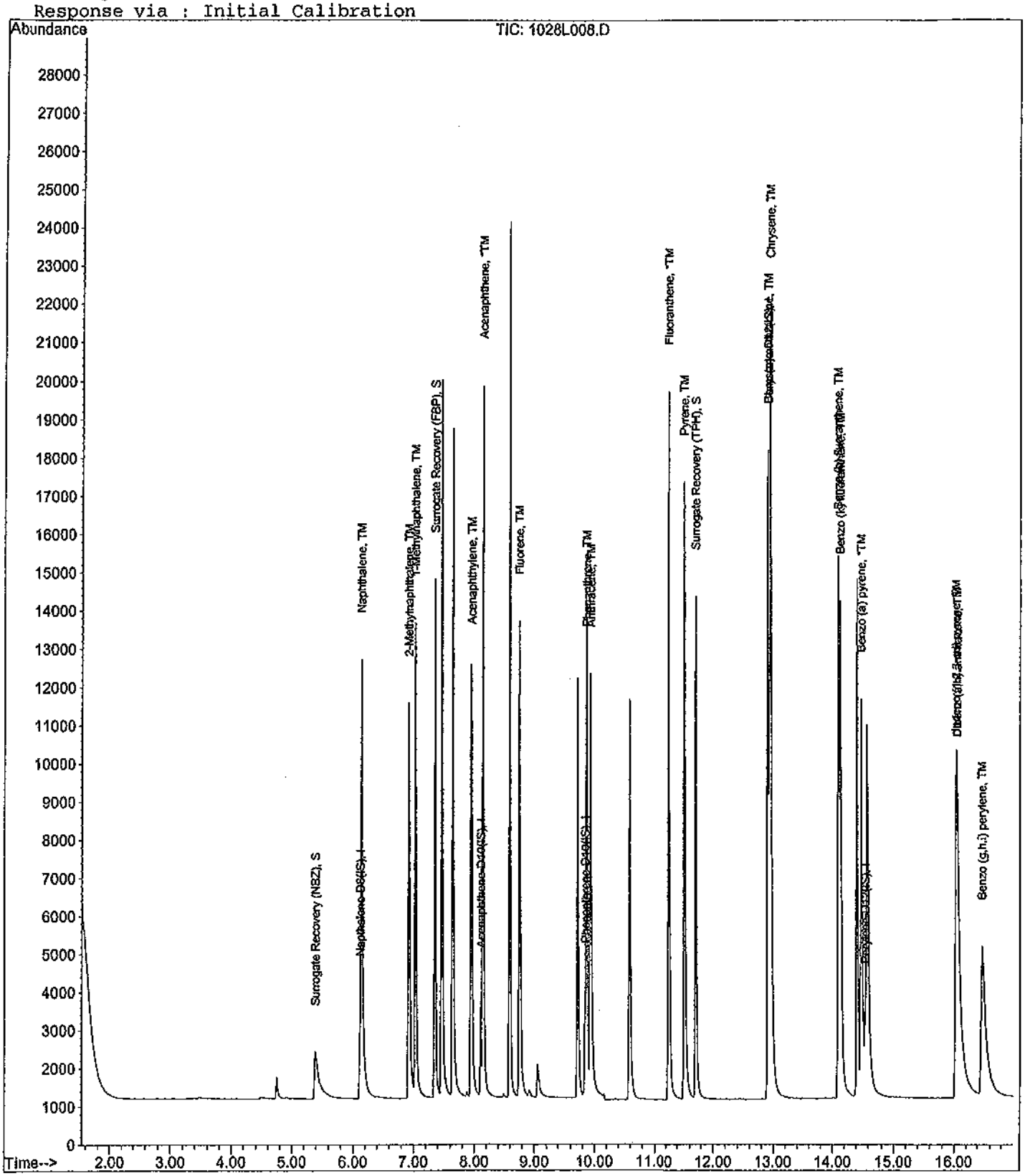
(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : M:\LINUS\DATA\L111027\1028L008.D Vial: 8
Acq On : 28 Oct 11 12:23 Operator: LF
Sample : 10ug/ml PAH Inst : Linus
Misc : Multiplr: 1.00

Quant Time: Oct 30 10:41 2011 Quant Results File: SIM2.RES

Method : M:\LINUS\DATA\L111027\SIM2.M (RTE Integrator)
Title : EPA 8270C
Last Update : Tue Nov 01 17:14:29 2011
Response via : Initial Calibration



Quantitation Report (QT Reviewed)

Data File : M:\LINUS\DATA\L111027\1028L009.D Vial: 9
Acq On : 28 Oct 11 12:49 Operator: LF
Sample : 50ug/ml PAH Inst : Linus
Misc : Multiplr: 1.00

Quant Time: Oct 30 10:41 2011 Quant Results File: SIM2.RES

Quant Method : M:\LINUS\DATA\L111028\SIM2.M (RTE Integrator)
Title : EPA 8270C
Last Update : Sun Oct 30 10:41:31 2011
Response via : Initial Calibration
DataAcq Meth : 87SIMAQ

Internal Standards	R.T.	QIon	Response	Conc Units		Dev(Min)
1) Napthalene-D8(IS)	6.11	136	2170	2.50000 ppb		-0.01
6) Acenaphthene-D10(IS)	8.11	164	955	2.50000 ppb		-0.02
11) Phenanthrene-D10(IS)	9.84	188	1764	2.50000 ppb		-0.04
15) Chrysene-D12(IS)	12.91	240	2325	2.50000 ppb		-0.05
21) Perylene-D12(IS)	14.54	264	1951	2.50000 ppb		-0.06
System Monitoring Compounds						
2) Surrogate Recovery (NBZ)	5.34	82	19569	80.30257 ppb		0.00
Spiked Amount 2.000			Recovery	= 4015.150%		
7) Surrogate Recovery (FBP)	7.34	172	37203	62.30259 ppb		-0.04
Spiked Amount 2.000			Recovery	= 3115.150%		
17) Surrogate Recovery (TPH)	11.70	244	43552	58.37048 ppb		-0.06
Spiked Amount 2.000			Recovery	= 2918.500%		
Target Compounds						Qvalue
3) Naphthalene	6.12	128	64981	36.91273 ppb		98
4) 2-Methylnaphthalene	6.92	142	39285	37.95912 ppb		91
5) 1-Methylnaphthalene	7.02	142	37731	34.00777 ppb		98
8) Acenaphthylene	7.94	152	59152	39.07406 ppb		100
9) Acenaphthene	8.13	154	32228	36.13782 ppb		90
10) Fluorene	8.75	166	36584	35.91740 ppb		95
12) Phenanthrene	9.86	178	48574	30.13920 ppb		99
13) Anthracene	9.92	178	49934	28.19038 ppb		99
14) Fluoranthene	11.23	202	84927	28.95874 ppb	#	86
16) Pyrene	11.50	202	87985	29.37950 ppb		93
18) Benz (a) anthracene	12.90	228	63776	37.84310 ppb		99
19) Chrysene	12.94	228	76944	35.03889 ppb		99
20) Indeno (1,2,3-cd) pyrene	16.01	276	67886	51.38427 ppb		97
22) Benzo (b) fluoranthene	14.09	252	68863	46.90706 ppb	#	96
23) Benzo (k) fluoranthene	14.12	252	60905	30.92236 ppb		100
24) Benzo (a) pyrene	14.45	252	61841	39.88811 ppb	#	94
25) Dibenz (a,h) anthracene	16.02	278	54590	55.19334 ppb		99
26) Benzo (g,h,i) perylene	16.44	276	56362	44.98303 ppb		98

(#) = qualifier out of range (m) = manual integration

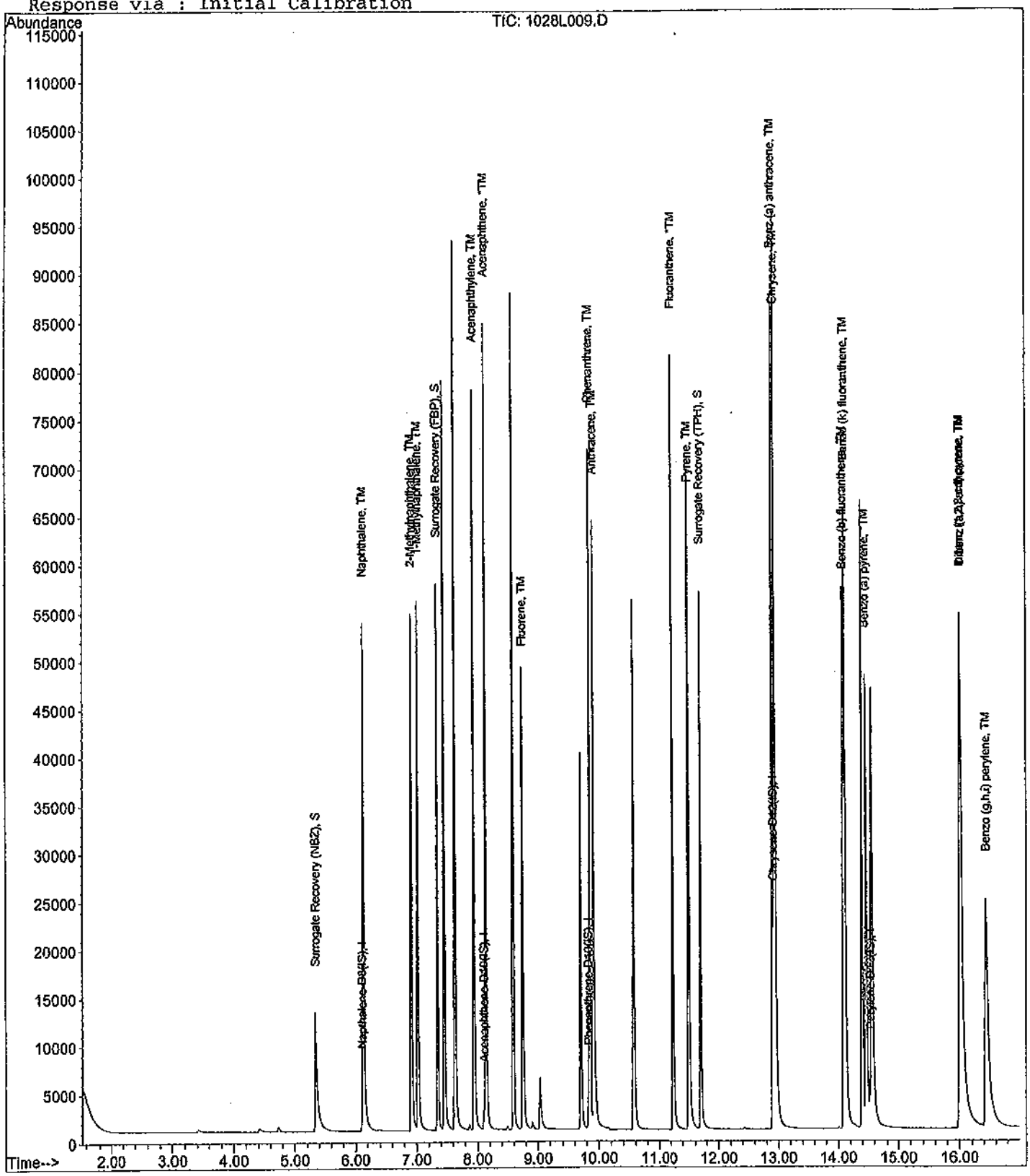
Quantitation Report

Data File : M:\LINUS\DATA\L111027\1028L009.D
Acq On : 28 Oct 11 12:49
Sample : 50ug/ml PAH
Misc :

Vial: 9
Operator: LF
Inst : Linus
Multiplr: 1.00

Quant Time: Oct 30 10:41 2011 Quant Results File: SIM2.RES

Method : M:\LINUS\DATA\L111027\SIM2.M (RTE Integrator)
Title : EPA 8270C
Last Update : Tue Nov 01 17:14:29 2011
Response via : Initial Calibration



Quantitation Report (QT Reviewed)

Data File : M:\LINUS\DATA\L111027\1028L010.D Vial: 10
Acq On : 28 Oct 11 13:14 Operator: LF
Sample : 100ug/ml PAH Inst : Linus
Misc : Multiplr: 1.00

Quant Time: Oct 30 10:42 2011 Quant Results File: SIM2.RES

Quant Method : M:\LINUS\DATA\L111028\SIM2.M (RTE Integrator)
Title : EPA 8270C
Last Update : Sun Oct 30 10:41:31 2011
Response via : Initial Calibration
DataAcq Meth : 87SIMAQ

Internal Standards	R.T.	QIon	Response	Conc Units		Dev(Min)
1) Napthalene-D8(IS)	6.11	136	2028	2.50000 ppb		-0.01
6) Acenaphthene-D10(IS)	8.11	164	919	2.50000 ppb		-0.02
11) Phenanthrene-D10(IS)	9.84	188	1786	2.50000 ppb		-0.04
15) Chrysene-D12(IS)	12.91	240	2218	2.50000 ppb		-0.05
21) Perylene-D12(IS)	14.54	264	1949	2.50000 ppb		-0.06
System Monitoring Compounds						
2) Surrogate Recovery (NBZ)	5.32	82	39811	174.48341 ppb		-0.01
Spiked Amount 2.000			Recovery	= 8724.150%		
7) Surrogate Recovery (FBP)	7.34	172	68503	119.21355 ppb		-0.04
Spiked Amount 2.000			Recovery	= 5960.700%		
17) Surrogate Recovery (TPH)	11.70	244	80239	112.72808 ppb		-0.06
Spiked Amount 2.000			Recovery	= 5636.400%		
Target Compounds						Qvalue
3) Naphthalene	6.12	128	118023	71.73782 ppb		98
4) 2-Methylnaphthalene	6.92	142	72350	74.80311 ppb		91
5) 1-Methylnaphthalene	7.02	142	67525	65.12327 ppb		99
8) Acenaphthylene	7.94	152	108807	74.69023 ppb		99
9) Acenaphthene	8.13	154	58631	68.31936 ppb		89
10) Fluorene	8.75	166	64716	66.02573 ppb		95
12) Phenanthrene	9.86	178	89156	54.63809 ppb		98
13) Anthracene	9.92	178	91266	50.88980 ppb		98
14) Fluoranthene	11.23	202	154470	52.02296 ppb	#	84
16) Pyrene	11.50	202	164055	57.42311 ppb	#	90
18) Benz (a) anthracene	12.90	228	140011	87.08694 ppb		99
19) Chrysene	12.94	228	127613	60.91607 ppb	#	95
20) Indeno (1,2,3-cd) pyrene	16.02	276	133093	105.60065 ppb	#	87
22) Benzo (b) fluoranthene	14.09	252	126697	86.39011 ppb		96
23) Benzo (k) fluoranthene	14.12	252	120651	61.31914 ppb	#	94
24) Benzo (a) pyrene	14.47	252	119503	77.15982 ppb		95
25) Dibenz (a,h) anthracene	16.03	278	107509	108.80876 ppb		91
26) Benzo (g,h,i) perylene	16.44	276	112699	90.03841 ppb		99

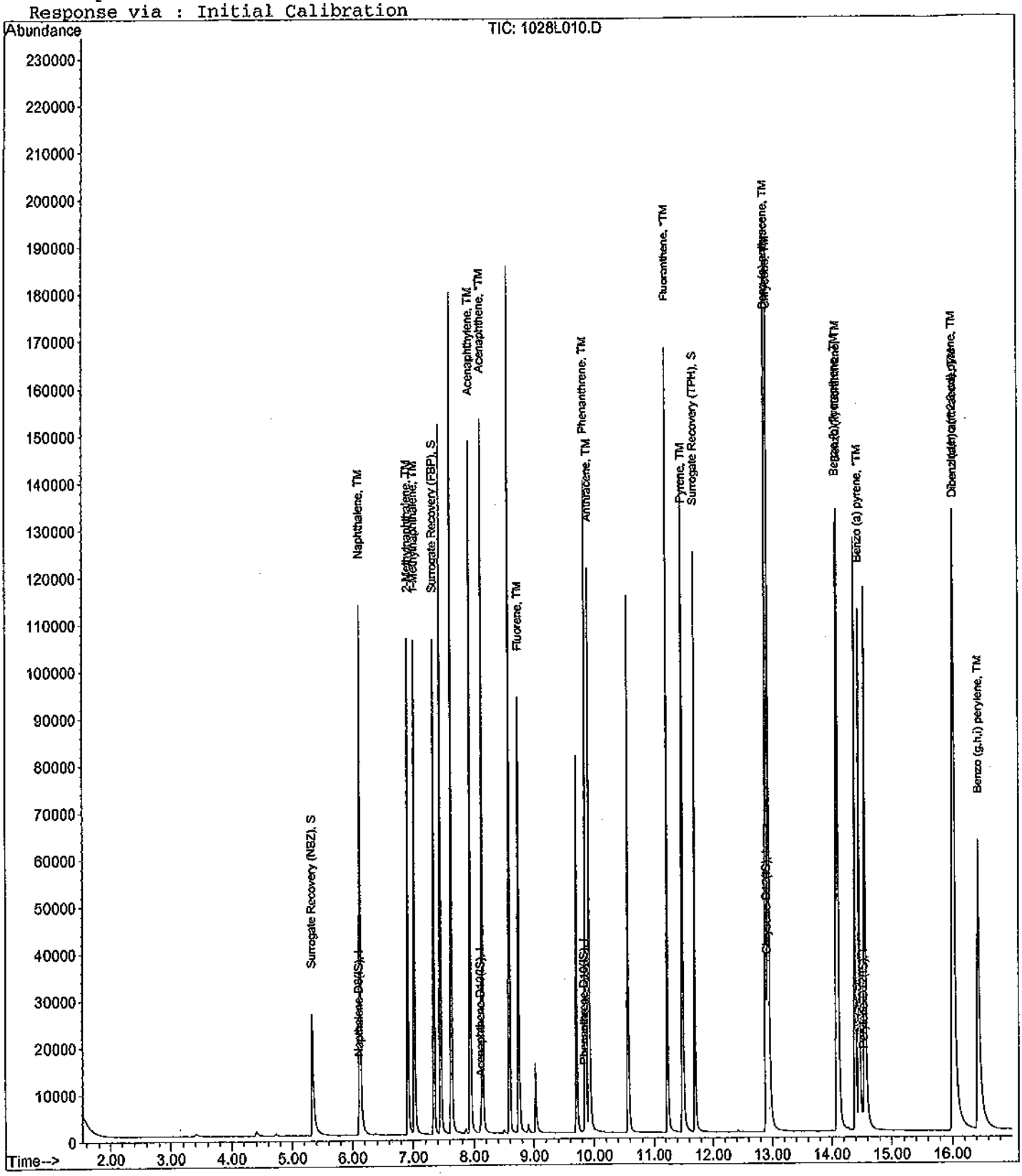
(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : M:\LINUS\DATA\L111027\1028L010.D Vial: 10
Acq On : 28 Oct 11 13:14 Operator: LF
Sample : 100ug/ml PAH Inst : Linus
Misc : Multiplr: 1.00

Quant Time: Oct 30 10:42 2011 Quant Results File: SIM2.RES

Method : M:\LINUS\DATA\L111027\SIM2.M (RTE Integrator)
Title : EPA 8270C
Last Update : Tue Nov 01 17:14:29 2011
Response via : Initial Calibration



Form 7

Second Source Calibration

Lab Name: APPL, Inc.
Case No:
Matrix:

SDG No: 66102
Date Analyzed: 10/28/11
Instrument: Linus
Initial Cal. Date: 10/27/11
Data File: 1028L011.D

		Compound	MEAN	CCRF	%D		%Drift
1	I	Napthalene-D8(IS)	ISTD			I	
2	TM	Naphthalene	1.742	1.546	11	TM	
3	TM	2-Methylnaphthalene	0.8931	0.8782	1.7	TM	
4	TM	1-Methylnaphthalene	1.031	1.007	2.4	TM	
5	I	Acenaphthene-D10(IS)	ISTD			I	
6	TM	Acenaphthylene	3.327	3.132	5.8	TM	
7	*TM	Acenaphthene	1.904	1.812	4.8	*TM	
8	TM	Fluorene	2.083	1.993	4.3	TM	
9	I	Phenanthrene-D10(IS)	ISTD			I	
10	TM	Phenanthrene	1.609	1.555	3.4	TM	
11	TM	Anthracene	1.634	1.624	0.64	TM	
12	*TM	Fluoranthene	2.792	2.916	4.4	*TM	
13	I	Chrysene-D12(IS)	ISTD			I	
14	TM	Pyrene	2.200	2.429	10	TM	
15	TM	Benz (a) anthracene	1.449	1.392	3.9	TM	
16	TM	Chrysene	1.939	2.190	13	TM	
17	TM	Indeno (1,2,3-cd) pyrene	1.502	1.468	2.3	TM	
18	I	Perylene-D12(IS)	ISTD			I	
19	TM	Benzo (b) fluoranthene	1.761	1.686	4.3	TM	
20	TM	Benzo (k) fluoranthene	1.823	2.176	19	TM	
21	*TM	Benzo (a) pyrene	1.723	1.689	1.9	*TM	
22	TM	Dibenz (a,h) anthracene	1.447	1.354	6.4	TM	
23	TM	Benzo (g,h,I) perylene	1.525	1.483	2.8	TM	
24							
25							
26							
27							
28							
29							
30							
31							
32							
33							
34							
35							
36							
37							
38							
39							
40							
		Average			5.7		

Quantitation Report (QT Reviewed)

Data File : M:\LINUS\DATA\L111027\1028L011.D Vial: 11
Acq On : 28 Oct 11 13:40 Operator: LF
Sample : 5.0ug/ml SS PAH 10-27-11 Inst : Linus
Misc : Multiplr: 1.00

Quant Time: Oct 30 11:17 2011 Quant Results File: SIM2.RES

Quant Method : M:\LINUS\DATA\L111027\SIM2.M (RTE Integrator)
Title : EPA 8270C
Last Update : Sun Oct 30 11:15:17 2011
Response via : Initial Calibration
DataAcq Meth : 87SIMAQ

Internal Standards	R.T.	QIon	Response	Conc Units		Dev(Min)
1) Napthalene-D8(IS)	6.13	136	2295	2.50000 ppb		0.01
6) Acenaphthene-D10(IS)	8.11	164	1033	2.50000 ppb		0.00
11) Phenanthrene-D10(IS)	9.85	188	1773	2.50000 ppb		0.00
15) Chrysene-D12(IS)	12.93	240	2205	2.50000 ppb		0.00
21) Perylene-D12(IS)	14.56	264	1840	2.50000 ppb		0.00
System Monitoring Compounds						
2) Surrogate Recovery (NBZ)	0.00	82	0d	0.00000 ppb		
Spiked Amount 2.000			Recovery =	0.000%		
7) Surrogate Recovery (FBP)	0.00	172	0d	0.00000 ppb		
Spiked Amount 2.000			Recovery =	0.000%		
17) Surrogate Recovery (TPH)	0.00	244	0d	0.00000 ppb		
Spiked Amount 2.000			Recovery =	0.000%		
Target Compounds						Qvalue
3) Naphthalene	6.14	128	7095	4.43732 ppb		99
4) 2-Methylnaphthalene	6.93	142	4031	4.91655 ppb		99
5) 1-Methylnaphthalene	7.04	142	4620	4.88168 ppb		94
8) Acenaphthylene	7.95	152	6471	4.70758 ppb		99
9) Acenaphthene	8.15	154	3744	4.75904 ppb		91
10) Fluorene	8.76	166	4117	4.78272 ppb		99
12) Phenanthrene	9.87	178	5514	4.83130 ppb		99
13) Anthracene	9.94	178	5757	4.96794 ppb		98
14) Fluoranthene	11.26	202	10339	5.22192 ppb		93
16) Pyrene	11.51	202	10711	5.51952 ppb	#	91
18) Benz (a) anthracene	12.93	228	6140	4.80346 ppb		99
19) Chrysene	12.96	228	9659	5.64891 ppb		99
20) Indeno (1,2,3-cd) pyrene	16.06	276	6475	4.88617 ppb	#	91
22) Benzo (b) fluoranthene	14.12	252	6204	4.78607 ppb		99
23) Benzo (k) fluoranthene	14.14	252	8006	5.96784 ppb	#	65
24) Benzo (a) pyrene	14.49	252	6217	4.90268 ppb		97
25) Dibenz (a,h) anthracene	16.08	278	4984	4.68078 ppb		96
26) Benzo (g,h,i) perylene	16.52	276	5458	4.86160 ppb		99

(#) = qualifier out of range (m) = manual integration

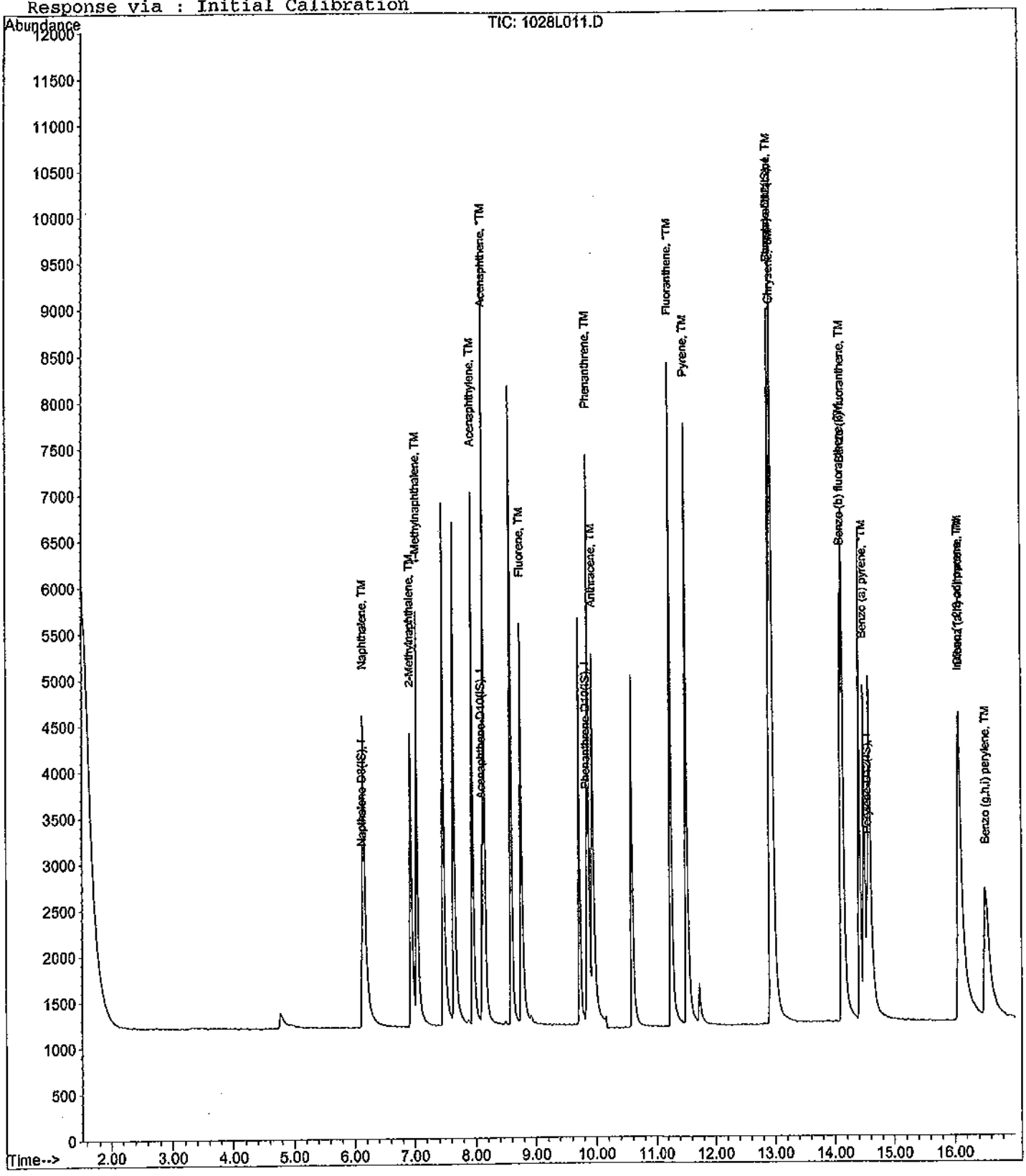
Quantitation Report

Data File : M:\LINUS\DATA\L111027\1028L011.D
Acq On : 28 Oct 11 13:40
Sample : 5.0ug/ml SS PAH 10-27-11
Misc :

Vial: 11
Operator: LF
Inst : Linus
Multiplr: 1.00

Quant Time: Oct 30 11:17 2011 Quant Results File: SIM2.RES

Method : M:\LINUS\DATA\L111027\SIM2.M (RTE Integrator)
Title : EPA 8270C
Last Update : Tue Nov 01 17:14:29 2011
Response via : Initial Calibration



Form 7

Continuing Calibration

Lab Name: APPL, Inc.
Case No:
Matrix:

SDG No: 66602
Date Analyzed: 5 Nov 11 16:54
Instrument: Linus
Initial Cal. Date: 10/27/11
Data File: 1105L020.D

		Compound	MEAN	CCRF	%D	%Drift	
1	I	Napthalene-D8(IS)	ISTD			I	
2	S	Surrogate Recovery (NBZ)	0.4477	0.4479	0.03	S	
3	TM	Naphthalene	1.742	1.518	13	TM	
4	TM	2-Methylnaphthalene	0.8931	0.9102	1.9	TM	
5	TM	1-Methylnaphthalene	1.031	0.9456	8.3	TM	
6	I	Acenaphthene-D10(IS)	ISTD			I	
7	S	Surrogate Recovery (FBP)	2.229	2.036	8.6	S	
8	TM	Acenaphthylene	3.327	3.080	7.4	TM	
9	*TM	Acenaphthene	1.904	1.695	11	*TM	
10	TM	Fluorene	2.083	1.961	5.9	TM	
11	I	Phenanthrene-D10(IS)	ISTD			I	
12	TM	Phenanthrene	1.609	1.494	7.2	TM	
13	TM	Anthracene	1.634	1.588	2.8	TM	
14	*TM	Fluoranthene	2.792	2.868	2.7	*TM	
15	I	Chrysene-D12(IS)	ISTD			I	
16	TM	Pyrene	2.200	2.193	0.31	TM	
17	S	Surrogate Recovery (TPH)	1.077	1.106	2.7	S	
18	TM	Benz (a) anthracene	1.449	1.737	20	TM	
19	TM	Chrysene	1.939	1.906	1.7	TM	
20	TM	Indeno (1,2,3-cd) pyrene	1.502	1.682	12	TM	
21	I	Perylene-D12(IS)	ISTD			I	
22	TM	Benzo (b) fluoranthene	1.761	1.921	9.1	TM	
23	TM	Benzo (k) fluoranthene	1.823	1.713	6.0	TM	
24	*TM	Benzo (a) pyrene	1.723	1.698	1.4	*TM	
25	TM	Dibenz (a,h) anthracene	1.447	1.482	2.5	TM	
26	TM	Benzo (g,h,i) perylene	1.525	1.535	0.66	TM	
27							
28							
29							
30							
31							
32							
33							
34							
35							
36							
37							
38							
39							
40							
		Average			6.0		

Quantitation Report (QT Reviewed)

Data File : M:\LINUS\DATA\L111027\1105L020.D Vial: 20
Acq On : 5 Nov 11 16:54 Operator: LF
Sample : 5.0ug/ml PAH 10-27-11 Inst : Linus
Misc : Multiplr: 1.00

Quant Time: Nov 8 9:47 2011 Quant Results File: SIM2.RES

Quant Method : M:\LINUS\DATA\L111027\SIM2.M (RTE Integrator)
Title : EPA 8270C
Last Update : Wed Nov 02 15:56:51 2011
Response via : Initial Calibration
DataAcq Meth : 87SIMAQ

Internal Standards	R.T.	QIon	Response	Conc Units	Dev(Min)
1) Napthalene-D8(IS)	6.11	136	1948	2.50000 ppb	-0.01
6) Acenaphthene-D10(IS)	8.11	164	877	2.50000 ppb	0.00
11) Phenanthrene-D10(IS)	9.85	188	1506	2.50000 ppb	0.00
15) Chrysene-D12(IS)	12.93	240	2029	2.50000 ppb	0.00
21) Perylene-D12(IS)	14.56	264	1837	2.50000 ppb	0.00
System Monitoring Compounds					
2) Surrogate Recovery (NBZ)	5.38	82	1745	5.00173 ppb	-0.06
Spiked Amount 2.000			Recovery	= 250.100%	
7) Surrogate Recovery (FBP)	7.35	172	3572	4.56887 ppb	0.00
Spiked Amount 2.000			Recovery	= 228.450%	
17) Surrogate Recovery (TPH)	11.70	244	4488	5.13686 ppb	-0.01
Spiked Amount 2.000			Recovery	= 256.850%	
Target Compounds					Qvalue
3) Naphthalene	6.13	128	5914	4.35756 ppb	99
4) 2-Methylnaphthalene	6.92	142	3546	5.09542 ppb	94
5) 1-Methylnaphthalene	7.02	142	3684	4.58607 ppb	92
8) Acenaphthylene	7.95	152	5403	4.62980 ppb	99
9) Acenaphthene	8.15	154	2973	4.45122 ppb	95
10) Fluorene	8.75	166	3440	4.70710 ppb	96
12) Phenanthrene	9.87	178	4499	4.64084 ppb	100
13) Anthracene	9.93	178	4783	4.85919 ppb	97
14) Fluoranthene	11.26	202	8637	5.13568 ppb	98
16) Pyrene	11.51	202	8901	4.98467 ppb	97
18) Benz (a) anthracene	12.91	228	7049	5.99294 ppb	98
19) Chrysene	12.96	228	7735	4.91608 ppb	98
20) Indeno (1,2,3-cd) pyrene	16.06	276	6825	5.59703 ppb #	91
22) Benzo (b) fluoranthene	14.10	252	7058	5.45378 ppb	97
23) Benzo (k) fluoranthene	14.14	252	6292	4.69785 ppb	95
24) Benzo (a) pyrene	14.49	252	6239	4.92807 ppb	99
25) Dibenz (a,h) anthracene	16.06	278	5446	5.12302 ppb	99
26) Benzo (g,h,i) perylene	16.51	276	5641	5.03281 ppb	96

(#) = qualifier out of range (m) = manual integration

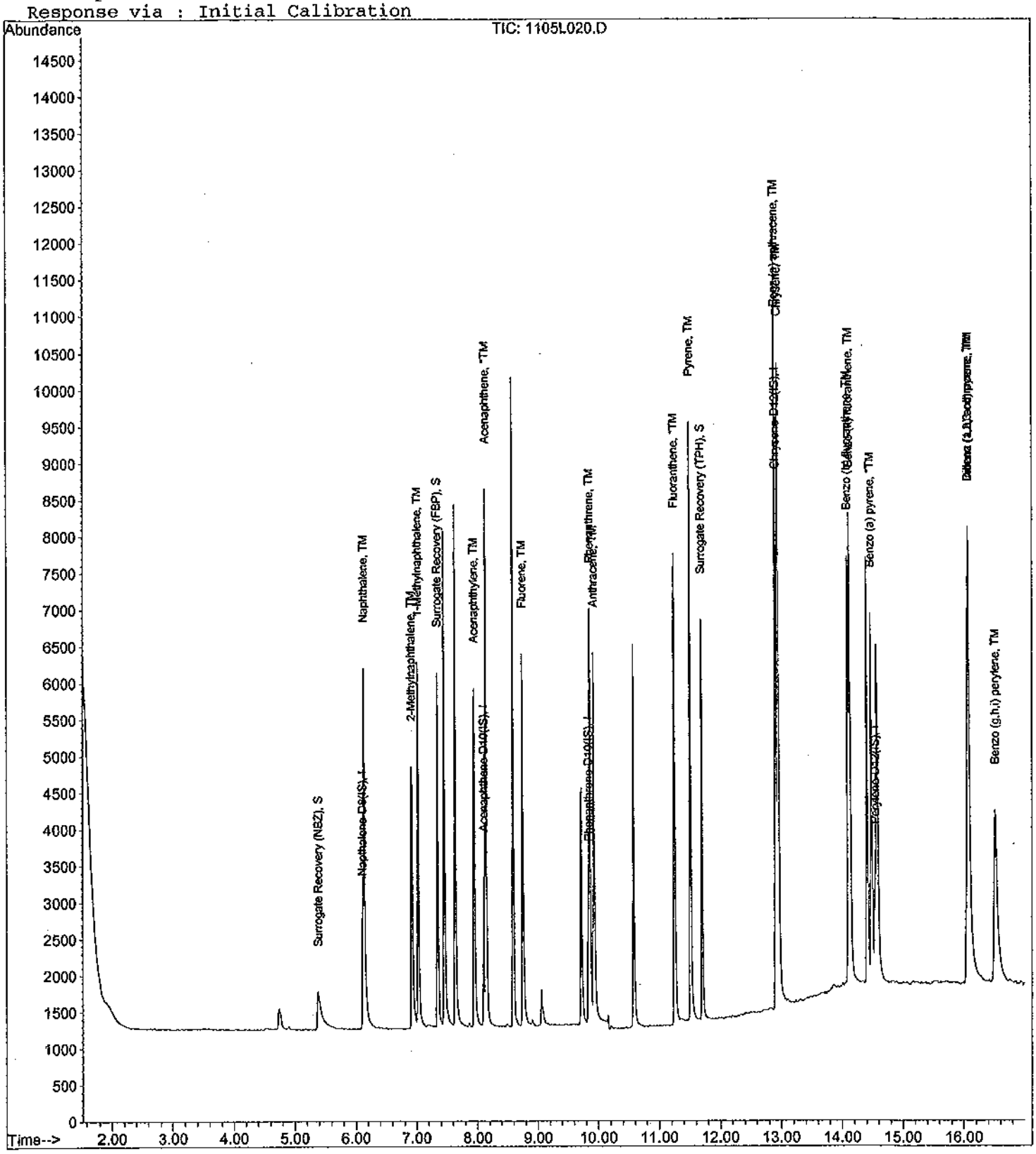
Quantitation Report

Data File : M:\LINUS\DATA\L111027\1105L020.D
Acq On : 5 Nov 11 16:54
Sample : 5.0ug/ml PAH 10-27-11
Misc :

Vial: 20
Operator: LF
Inst : Linus
Multiplr: 1.00

Quant Time: Nov 8 9:47 2011 Quant Results File: SIM2.RES

Method : M:\LINUS\DATA\L111027\SIM2.M (RTE Integrator)
Title : EPA 8270C
Last Update : Tue Nov 08 16:22:04 2011
Response via : Initial Calibration



EPA METHOD 8270
Polynuclear Aromatic Hydrocarbons
Raw Data

APPL, INC.

Method Blank
EPA 8270D SIM

APPL Inc.
908 North Temperance Avenu
Clovis, CA 93611

Blank Name/QCG: **111031W-49334 - 161019**
Batch ID: #SIMHC-111031A

Sample Type	Analyte	Result	LOQ	LOD	DL	Units	Extraction Date	Analysis Date
BLANK	1-METHYLNAPHTHALENE	0.12 U	0.2	0.12	0.06	ug/L	10/31/11	11/05/11
BLANK	2-METHYLNAPHTHALENE	0.12 U	0.2	0.12	0.06	ug/L	10/31/11	11/05/11
BLANK	ACENAPHTHENE	0.12 U	0.2	0.12	0.06	ug/L	10/31/11	11/05/11
BLANK	ACENAPHTHYLENE	0.12 U	0.2	0.12	0.06	ug/L	10/31/11	11/05/11
BLANK	ANTHRACENE	0.10 U	0.2	0.10	0.05	ug/L	10/31/11	11/05/11
BLANK	BENZO(A)ANTHRACENE	0.14 U	0.2	0.14	0.07	ug/L	10/31/11	11/05/11
BLANK	BENZO(A)PYRENE	0.14 U	0.2	0.14	0.07	ug/L	10/31/11	11/05/11
BLANK	BENZO(B)FLUORANTHENE	0.12 U	0.2	0.12	0.06	ug/L	10/31/11	11/05/11
BLANK	BENZO(GHI)PERYLENE	0.16 U	0.2	0.16	0.08	ug/L	10/31/11	11/05/11
BLANK	BENZO(K)FLUORANTHENE	0.14 U	0.2	0.14	0.07	ug/L	10/31/11	11/05/11
BLANK	CHRYSENE	0.10 U	0.2	0.10	0.05	ug/L	10/31/11	11/05/11
BLANK	DIBENZ(A,H)ANTHRACENE	0.10 U	0.2	0.10	0.05	ug/L	10/31/11	11/05/11
BLANK	FLUORANTHENE	0.16 U	0.2	0.16	0.08	ug/L	10/31/11	11/05/11
BLANK	FLUORENE	0.12 U	0.2	0.12	0.06	ug/L	10/31/11	11/05/11
BLANK	INDENO(1,2,3-CD)PYRENE	0.14 U	0.2	0.14	0.07	ug/L	10/31/11	11/05/11
BLANK	NAPHTHALENE	0.10 U	0.2	0.10	0.05	ug/L	10/31/11	11/05/11
BLANK	PHENANTHRENE	0.14 U	0.2	0.14	0.07	ug/L	10/31/11	11/05/11
BLANK	PYRENE	0.16 U	0.2	0.16	0.08	ug/L	10/31/11	11/05/11
BLANK	SURROGATE: 2-FLUORBIPHENY	51.7	50-110			%	10/31/11	11/05/11
BLANK	SURROGATE: NITROBENZENE-	66.3	40-110			%	10/31/11	11/05/11
BLANK	SURROGATE: TERPHENYL-D14 (	54.5	50-135			%	10/31/11	11/05/11

Quant Method:SIM2.M
Run #:1105L028
Instrument:Linus
Sequence:L111027
Initials:LF

GC SC-Blank-REG MDLs
Printed: 11/09/11 3:59:20 PM

Quantitation Report (QT Reviewed)

Data File : M:\LINUS\DATA\L111027\1105L028.D
Acq On : 5 Nov 11 20:15
Sample : 111031A BLK 1/1000
Misc :

Vial: 28
Operator: LF
Inst : Linus
Multiplr: 1.00

Quant Time: Nov 9 8:48 2011 Quant Results File: SIM2.RES

Quant Method : M:\LINUS\DATA\L111027\SIM2.M (RTE Integrator)
Title : EPA 8270C
Last Update : Wed Nov 02 15:56:51 2011
Response via : Initial Calibration
DataAcq Meth : 87SIMAQ

Internal Standards	R.T.	QIon	Response	Conc Units	Dev(Min)
1) Napthalene-D8(IS)	6.12	136	2305	2.50000 ppb	0.00
6) Acenaphthene-D10(IS)	8.11	164	1068	2.50000 ppb	0.00
11) Phenanthrene-D10(IS)	9.86	188	2122	2.50000 ppb	0.01
15) Chrysene-D12(IS)	12.94	240	2454	2.50000 ppb	0.01
21) Perylene-D12(IS)	14.57	264	2143	2.50000 ppb	0.01
System Monitoring Compounds					
2) Surrogate Recovery (NBZ)	5.42	82	547	1.32504 ppb	-0.02
Spiked Amount 2.000			Recovery = 66.250%		
7) Surrogate Recovery (FBP)	7.36	172	984	1.03352 ppb	0.01
Spiked Amount 2.000			Recovery = 51.700%		
17) Surrogate Recovery (TPH)	11.71	244	1151	1.08925 ppb	0.00
Spiked Amount 2.000			Recovery = 54.450%		

Target Compounds Qvalue

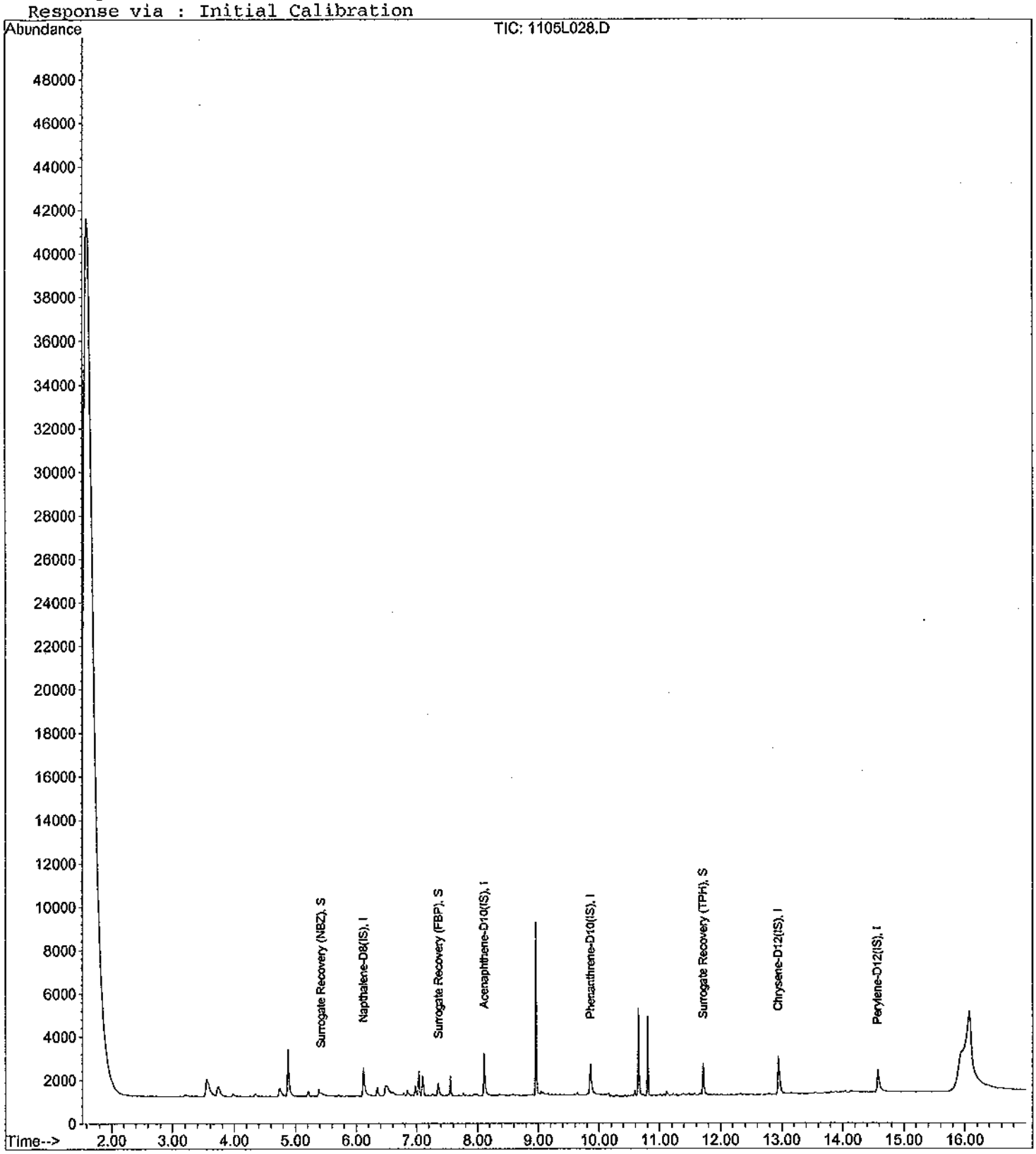
(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : M:\LINUS\DATA\L111027\1105L028.D Vial: 28
Acq On : 5 Nov 11 20:15 Operator: LF
Sample : 111031A BLK 1/1000 Inst : Linus
Misc : Multiplr: 1.00

Quant Time: Nov 9 8:48 2011 Quant Results File: SIM2.RES

Method : M:\LINUS\DATA\L111027\SIM2.M (RTE Integrator)
Title : EPA 8270C
Last Update : Tue Nov 08 16:22:04 2011
Response via : Initial Calibration



Laboratory Control Spike Recovery
EPA 8270D SIM

APPL ID: **111031W-49334 LCS - 161019**
Batch ID: #SIMHC-111031A

APPL Inc.
908 North Temperance Avenue
Clovis, CA 93611

Compound Name	Spike Level ug/L	SPK Result ug/L	SPK % Recovery	Recovery Limits
1-METHYLNAPHTHALENE	4.00	2.25	56.3	45-105
2-METHYLNAPHTHALENE	4.00	2.26	56.5	45-105
ACENAPHTHENE	4.00	2.55	63.7	45-110
ACENAPHTHYLENE	4.00	2.39	59.8	50-105
ANTHRACENE	4.00	2.47	61.8	55-110
BENZO(A)ANTHRACENE	4.00	2.74	68.5	55-110
BENZO(A)PYRENE	4.00	2.48	62.0	55-110
BENZO(B)FLUORANTHENE	4.00	2.43	60.8	45-120
BENZO(GHI)PERYLENE	4.00	2.80	70.0	40-125
BENZO(K)FLUORANTHENE	4.00	3.23	80.8	45-125
CHRYSENE	4.00	2.86	71.5	55-110
DIBENZ(A,H)ANTHRACENE	4.00	2.89	72.3	40-125
FLUORANTHENE	4.00	2.86	71.5	55-115
FLUORENE	4.00	2.59	64.8	50-110
INDENO(1,2,3-CD)PYRENE	4.00	2.95	73.8	45-125
NAPHTHALENE	4.00	2.30	57.5	40-100
PHENANTHRENE	4.00	2.43	60.8	50-115
PYRENE	4.00	2.57	64.3	50-130
SURROGATE: 2-FLUORBIPHENYL (S)	2.00	1.11	55.5	50-110
SURROGATE: NITROBENZENE-D5 (S)	2.00	1.14	57.0	40-110
SURROGATE: TERPHENYL-D14 (S)	2.00	1.06	53.0	50-135

Comments: _____

Primary	SPK
Quant Method :	SIM2.M
Extraction Date :	10/31/11
Analysis Date :	11/05/11
Instrument :	Linus
Run :	1105L029
Initials :	LF

Printed: 11/09/11 3:59:21 PM

APPL Standard LCS

Quantitation Report (QT Reviewed)

Data File : M:\LINUS\DATA\L111027\1105L029.D Vial: 29
Acq On : 5 Nov 11 20:41 Operator: LF
Sample : 111031A LCS-1 1/1000 Inst : Linus
Misc : Multiplr: 1.00

Quant Time: Nov 9 8:50 2011 Quant Results File: SIM2.RES

Quant Method : M:\LINUS\DATA\L111027\SIM2.M (RTE Integrator)
Title : EPA 8270C
Last Update : Wed Nov 02 15:56:51 2011
Response via : Initial Calibration
DataAcq Meth : 87SIMAQ

Internal Standards	R.T.	QIon	Response	Conc	Units		Dev(Min)
1) Napthalene-D8(IS)	6.12	136	2079	2.50000	ppb		0.00
6) Acenaphthene-D10(IS)	8.11	164	961	2.50000	ppb		0.00
11) Phenanthrene-D10(IS)	9.85	188	1713	2.50000	ppb		0.00
15) Chrysene-D12(IS)	12.93	240	2367	2.50000	ppb		0.00
21) Perylene-D12(IS)	14.56	264	2017	2.50000	ppb		0.00
System Monitoring Compounds							
2) Surrogate Recovery (NBZ)	5.43	82	426	1.14411	ppb		-0.01
Spiked Amount 2.000			Recovery = 57.200%				
7) Surrogate Recovery (FBP)	7.36	172	951	1.11008	ppb		0.01
Spiked Amount 2.000			Recovery = 55.500%				
17) Surrogate Recovery (TPH)	11.71	244	1078	1.05766	ppb		0.00
Spiked Amount 2.000			Recovery = 52.900%				
Target Compounds							Qvalue
3) Naphthalene	6.14	128	3335	2.30246	ppb		99
4) 2-Methylnaphthalene	6.93	142	1681	2.26331	ppb		89
5) 1-Methylnaphthalene	7.04	142	1929	2.25003	ppb		88
8) Acenaphthylene	7.95	152	3060	2.39290	ppb		99
9) Acenaphthene	8.15	154	1866	2.54960	ppb		93
10) Fluorene	8.76	166	2077	2.59363	ppb		97
12) Phenanthrene	9.87	178	2685	2.43497	ppb		98
13) Anthracene	9.94	178	2766	2.47049	ppb		99
14) Fluoranthene	11.26	202	5473	2.86107	ppb		94
16) Pyrene	11.51	202	5362	2.57400	ppb	#	87
18) Benz (a) anthracene	12.93	228	3766	2.74459	ppb		98
19) Chrysene	12.96	228	5250	2.86024	ppb		99
20) Indeno (1,2,3-cd) pyrene	16.09	276	4203	2.95460	ppb	#	95
22) Benzo (b) fluoranthene	14.12	252	3457	2.43287	ppb	#	91
23) Benzo (k) fluoranthene	14.15	252	4746	3.22732	ppb		95
24) Benzo (a) pyrene	14.50	252	3444	2.47758	ppb		98
25) Dibenz (a,h) anthracene	16.09	278	3371	2.88809	ppb		95
26) Benzo (g,h,i) perylene	16.53	276	3449	2.80253	ppb		98

$\frac{3335 \times 2.5}{2079 \times 1.742} = 2.30$ VF 11/9/11

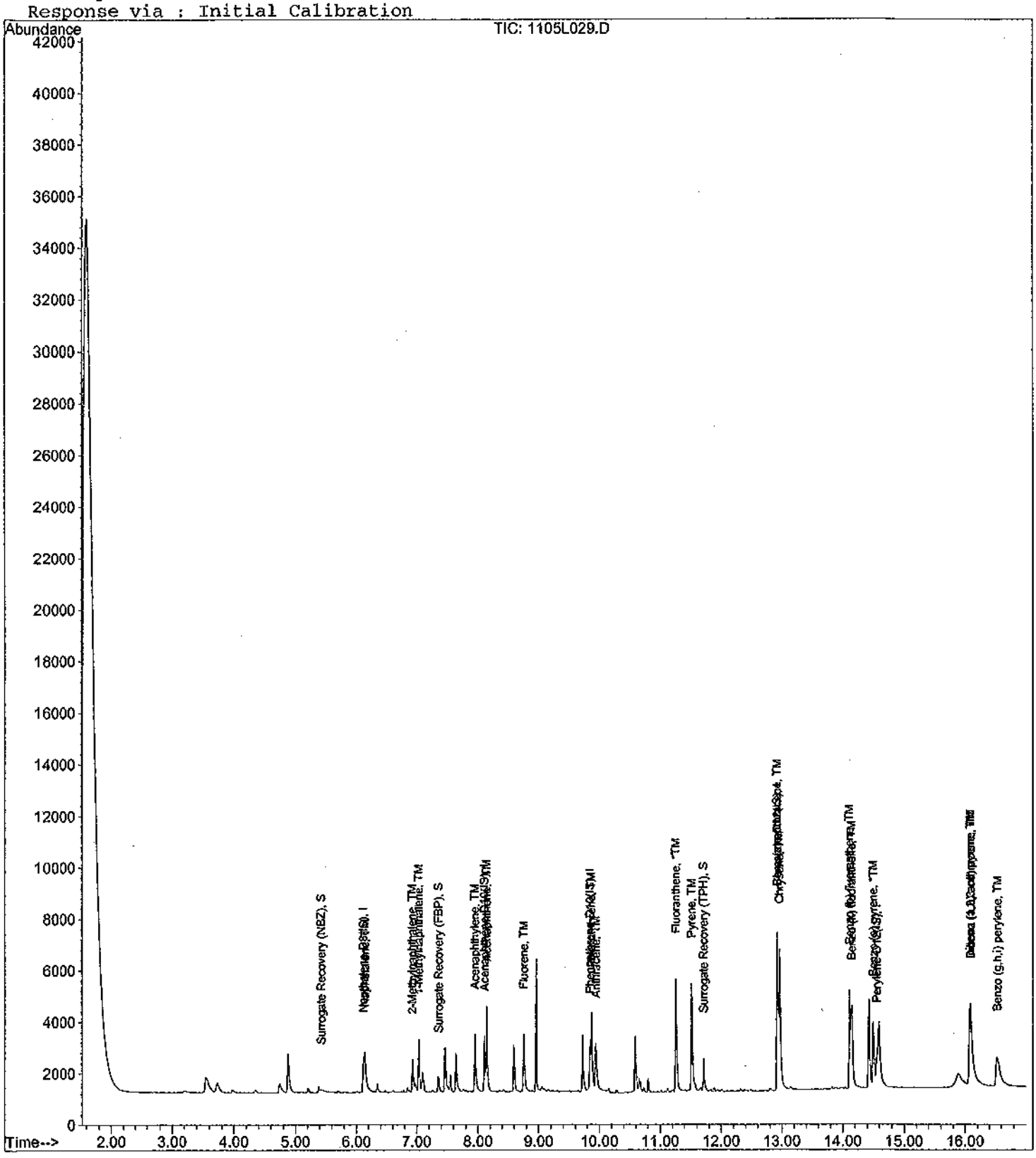
(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : M:\LINUS\DATA\L111027\1105L029.D Vial: 29
Acq On : 5 Nov 11 20:41 Operator: LF
Sample : 111031A LCS-1 1/1000 Inst : Linus
Misc : Multiplr: 1.00

Quant Time: Nov 9 8:50 2011 Quant Results File: SIM2.RES

Method : M:\LINUS\DATA\L111027\SIM2.M (RTE Integrator)
Title : EPA 8270C
Last Update : Tue Nov 08 16:22:04 2011
Response via : Initial Calibration



Matrix Spike Recoveries

EPA 8270D SIM

APPL ID: **111031W-49334 MS - 161019**
Batch ID: #SIMHC-111031A
Sample ID: AY49334
Client ID: ES047

APPL Inc.
908 North Temperance Avenue
Clovis, CA 93611

Compound Name	Spike Lvl ug/L	Matrix Result ug/L	SPK Result ug/L	DUP Result ug/L	SPK % Recovery	DUP % Recovery	Recovery Limits	RPD %	RPD Limits
1-METHYLNAPHTHALENE	3.88	0.62	2.88	2.71	58.2	53.9	45-105	6.1	25
2-METHYLNAPHTHALENE	3.88	0.20	2.26	1.98	53.1	45.9	45-105	13.2	25
ACENAPHTHENE	3.88	0.16	2.17	1.79	51.8	42.0 #	45-110	19.2	25
ACENAPHTHYLENE	3.88	ND	2.14	1.77	55.2	45.6 #	50-105	18.9	25
ANTHRACENE	3.88	ND	2.14	1.96	55.2	50.5 #	55-110	8.8	25
BENZO(A)ANTHRACENE	3.88	ND	3.04	2.84	78.4	73.2	55-110	6.8	25
BENZO(A)PYRENE	3.88	ND	2.62	2.30	67.5	59.3	55-110	13.0	25
BENZO(B)FLUORANTHENE	3.88	ND	2.72	2.34	70.1	60.3	45-120	15.0	25
BENZO(GHI)PERYLENE	3.88	ND	2.84	2.54	73.2	65.5	40-125	11.2	25
BENZO(K)FLUORANTHENE	3.88	ND	3.07	2.84	79.1	73.2	45-125	7.8	25
CHRYSENE	3.88	ND	2.75	2.56	70.9	66.0	55-110	7.2	25
DIBENZ(A,H)ANTHRACENE	3.88	ND	2.90	2.51	74.7	64.7	40-125	14.4	25
FLUORANTHENE	3.88	ND	2.84	2.59	73.2	66.8	55-115	9.2	25
FLUORENE	3.88	0.083	2.42	2.08	60.2	51.5	50-110	15.1	25
INDENO(1,2,3-CD)PYRENE	3.88	ND	2.93	2.71	75.5	69.8	45-125	7.8	25
NAPHTHALENE	3.88	1.0	3.70	2.56	69.6	40.2	40-100	36.4 #	25
PHENANTHRENE	3.88	ND	2.38	2.20	61.3	56.7	50-115	7.9	25
PYRENE	3.88	ND	2.63	2.45	67.8	63.1	50-130	7.1	25
SURROGATE: 2-FLUORBIPHENYL (S)	1.94	NA	1.04	1.08	53.6	55.7	50-110		
SURROGATE: NITROBENZENE-D5 (S)	1.94	NA	1.15	1.03	59.3	53.1	40-110		
SURROGATE: TERPHENYL-D14 (S)	1.94	NA	0.984	1.20	50.7	61.9	50-135		

= Recovery is outside QC limits.

Comments:

Primary	SPK	DUP
Quant Method :	SIM2.M	SIM2.M
Extraction Date :	10/31/11	10/31/11
Analysis Date :	11/05/11	11/06/11
Instrument :	Linus	Linus
Run :	1105L036	1105L037
Initials :	LF	

Printed: 11/09/11 4:01:35 PM
APPL MSD SCII

Quantitation Report (QT Reviewed)

Data File : M:\LINUS\DATA\L111027\1105L036.D Vial: 36
Acq On : 5 Nov 11 23:36 Operator: LF
Sample : AY49334W30 MS-1 1/1030 Inst : Linus
Misc : Multiplr: 0.97

Quant Time: Nov 9 9:18 2011 Quant Results File: SIM2.RES

Quant Method : M:\LINUS\DATA\L111027\SIM2.M (RTE Integrator)
Title : EPA 8270C
Last Update : Wed Nov 02 15:56:51 2011
Response via : Initial Calibration
DataAcq Meth : 87SIMAQ

Internal Standards	R.T.	QIon	Response	Conc Units		Dev(Min)
1) Napthalene-D8(IS)	6.11	136	2236	2.50000 ppb		-0.01
6) Acenaphthene-D10(IS)	8.11	164	1033	2.50000 ppb		0.00
11) Phenanthrene-D10(IS)	9.84	188	1791	2.50000 ppb		-0.01
15) Chrysene-D12(IS)	12.93	240	2429	2.50000 ppb		0.00
21) Perylene-D12(IS)	14.56	264	2046	2.50000 ppb		0.00
System Monitoring Compounds						
2) Surrogate Recovery (NBZ)	5.41	82	474	1.14916 ppb		-0.04
Spiked Amount 1.942			Recovery = 59.174%			
7) Surrogate Recovery (FBP)	7.35	172	984	1.03741 ppb		0.00
Spiked Amount 1.942			Recovery = 53.406%			
17) Surrogate Recovery (TPH)	11.70	244	1060	0.98394 ppb		-0.01
Spiked Amount 1.942			Recovery = 50.676%			
Target Compounds						Qvalue
3) Naphthalene	6.13	128	5940	3.70192 ppb	#	90
4) 2-Methylnaphthalene	6.92	142	1859	2.25943 ppb		91
5) 1-Methylnaphthalene	7.02	142	2737	2.88187 ppb		92
8) Acenaphthylene	7.94	152	3027	2.13796 ppb		95
9) Acenaphthene	8.15	154	1758	2.16952 ppb		99
10) Fluorene	8.75	166	2145	2.41926 ppb		98
12) Phenanthrene	9.86	178	2830	2.38319 ppb		96
13) Anthracene	9.93	178	2585	2.14395 ppb		99
14) Fluoranthene	11.25	202	5849	2.83927 ppb	#	80
16) Pyrene	11.51	202	5786	2.62780 ppb		95
18) Benz (a) anthracene	12.91	228	4407	3.03859 ppb		97
19) Chrysene	12.96	228	5335	2.74985 ppb		98
20) Indeno (1,2,3-cd) pyrene	16.08	276	4399	2.92567 ppb	#	93
22) Benzo (b) fluoranthene	14.12	252	4037	2.71919 ppb	#	91
23) Benzo (k) fluoranthene	14.14	252	4718	3.07067 ppb		96
24) Benzo (a) pyrene	14.50	252	3802	2.61781 ppb		97
25) Dibenz (a,h) anthracene	16.08	278	3537	2.90034 ppb		98
26) Benzo (g,h,i) perylene	16.53	276	3653	2.84098 ppb		99

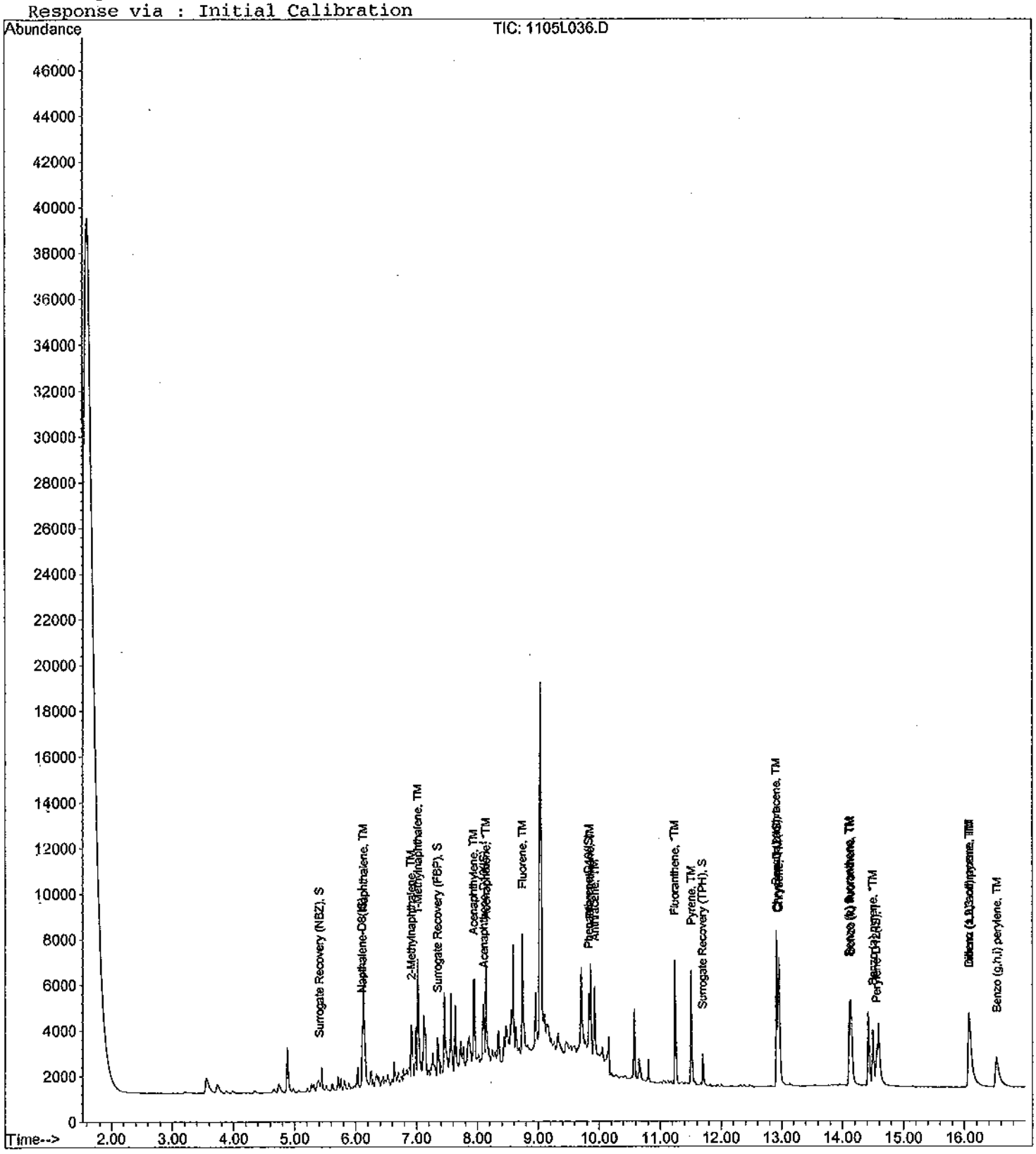
(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : M:\LINUS\DATA\L111027\1105L036.D Vial: 36
Acq On : 5 Nov 11 23:36 Operator: LF
Sample : AY49334W30 MS-1 1/1030 Inst : Linus
Misc : Multiplr: 0.97

Quant Time: Nov 9 9:18 2011 Quant Results File: SIM2.RES

Method : M:\LINUS\DATA\L111027\SIM2.M (RTE Integrator)
Title : EPA 8270C
Last Update : Tue Nov 08 16:22:04 2011
Response via : Initial Calibration



Quantitation Report (QT Reviewed)

Data File : M:\LINUS\DATA\L111027\1105L037.D Vial: 37
Acq On : 6 Nov 11 00:01 Operator: LF
Sample : AY49334W34 MSD-1 1/1030 Inst : Linus
Misc : Multiplr: 0.97

Quant Time: Nov 9 9:19 2011 Quant Results File: SIM2.RES

Quant Method : M:\LINUS\DATA\L111027\SIM2.M (RTE Integrator)
Title : EPA 8270C
Last Update : Wed Nov 02 15:56:51 2011
Response via : Initial Calibration
DataAcq Meth : 87SIMAQ

Internal Standards	R.T.	QIon	Response	Conc Units	Dev(Min)
1) Napthalene-D8(IS)	6.11	136	2287	2.50000 ppb	-0.01
6) Acenaphthene-D10(IS)	8.11	164	1131	2.50000 ppb	0.00
11) Phenanthrene-D10(IS)	9.83	188	1879	2.50000 ppb	-0.01
15) Chrysene-D12(IS)	12.93	240	2428	2.50000 ppb	0.00
21) Perylene-D12(IS)	14.56	264	2160	2.50000 ppb	0.00
System Monitoring Compounds					
2) Surrogate Recovery (NBZ)	5.41	82	433	1.02635 ppb	-0.04
Spiked Amount 1.942			Recovery = 52.839%		
7) Surrogate Recovery (FBP)	7.35	172	1118	1.07656 ppb	0.00
Spiked Amount 1.942			Recovery = 55.466%		
17) Surrogate Recovery (TPH)	11.70	244	1293	1.20071 ppb	-0.01
Spiked Amount 1.942			Recovery = 61.852%		
Target Compounds					Qvalue
3) Naphthalene	6.13	128	4199	2.55854 ppb #	85
4) 2-Methylnaphthalene	6.92	142	1664	1.97733 ppb	92
5) 1-Methylnaphthalene	7.02	142	2628	2.70539 ppb	91
8) Acenaphthylene	7.94	152	2746	1.77144 ppb	94
9) Acenaphthene	8.15	154	1590	1.79217 ppb	99
10) Fluorene	8.75	166	2022	2.08293 ppb	98
12) Phenanthrene	9.86	178	2744	2.20254 ppb	96
13) Anthracene	9.93	178	2473	1.95500 ppb	99
14) Fluoranthene	11.25	202	5590	2.58646 ppb #	83
16) Pyrene	11.51	202	5394	2.45077 ppb	94
18) Benz (a) anthracene	12.91	228	4123	2.84394 ppb	98
19) Chrysene	12.96	228	4973	2.56432 ppb	99
20) Indeno (1,2,3-cd) pyrene	16.08	276	4079	2.71396 ppb #	90
22) Benzo (b) fluoranthene	14.12	252	3669	2.34088 ppb #	90
23) Benzo (k) fluoranthene	14.14	252	4610	2.84203 ppb	96
24) Benzo (a) pyrene	14.50	252	3531	2.30290 ppb	98
25) Dibenz (a,h) anthracene	16.08	278	3235	2.51269 ppb	99
26) Benzo (g,h,i) perylene	16.53	276	3447	2.53929 ppb	97

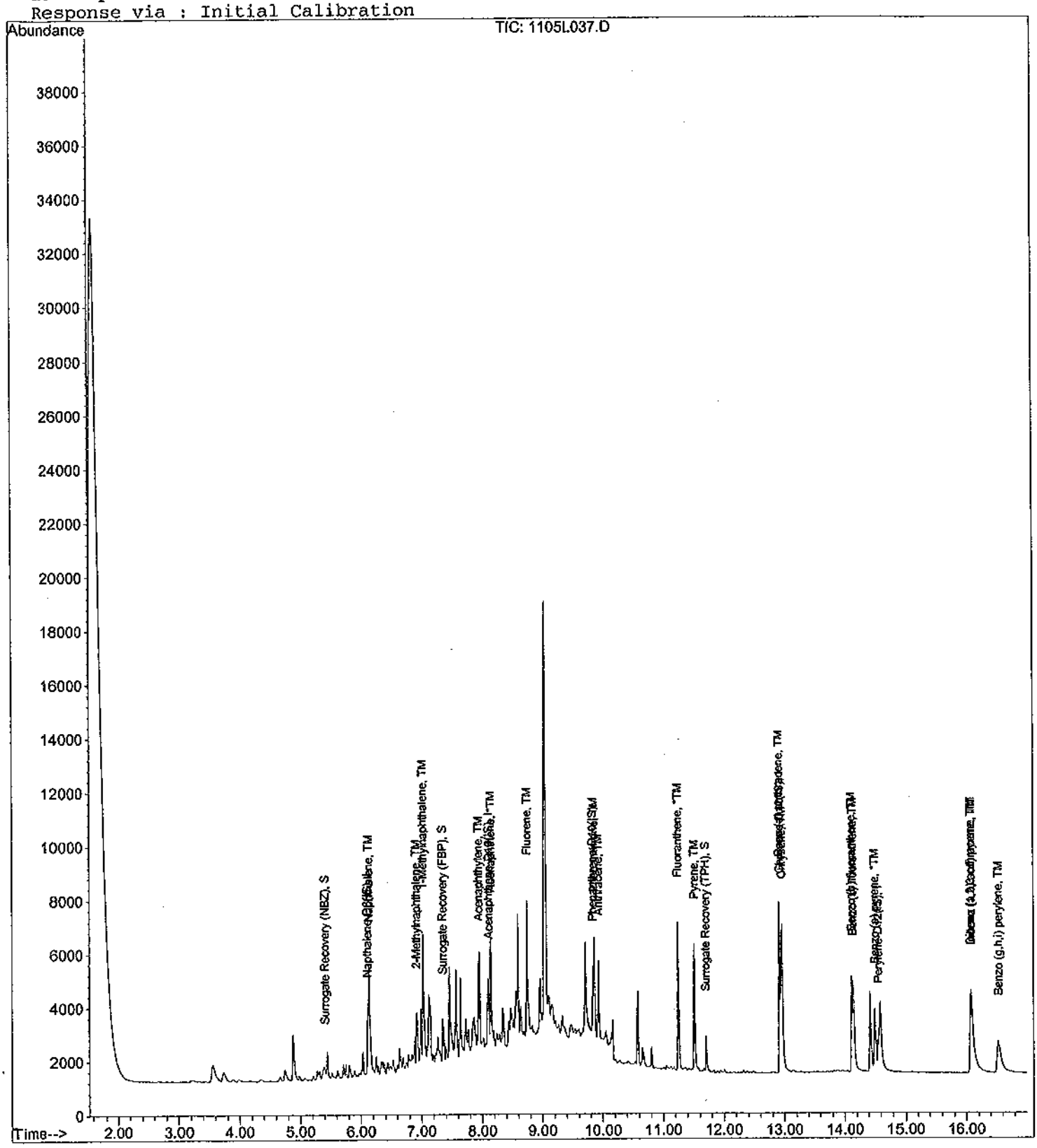
(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : M:\LINUS\DATA\L111027\1105L037.D Vial: 37
Acq On : 6 Nov 11 00:01 Operator: LF
Sample : AY49334W34 MSD-1 1/1030 Inst : Linus
Misc : Multiplr: 0.97

Quant Time: Nov 9 9:19 2011 Quant Results File: SIM2.RES

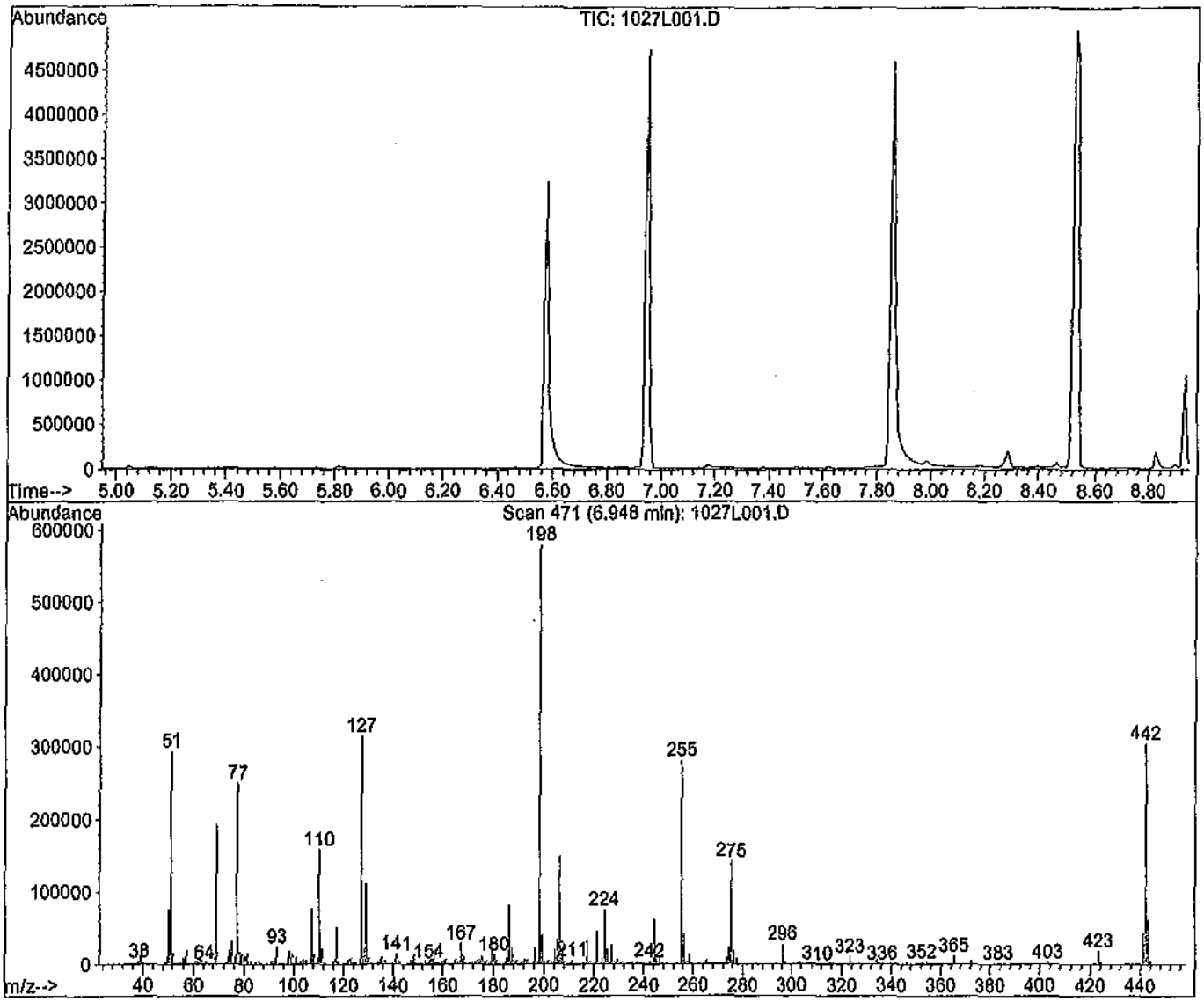
Method : M:\LINUS\DATA\L111027\SIM2.M (RTE Integrator)
Title : EPA 8270C
Last Update : Tue Nov 08 16:22:04 2011
Response via : Initial Calibration



DFTPP

Data File : M:\LINUS\DATA\L111027\1027L001.D Vial: 1
Acq On : 27 Oct 11 18:29 Operator: LF
Sample : SVTUNE 10-27-11 Inst : Linus
Misc : Multiplr: 1.00

Method : M:\LINUS\DATA\L111027\SIM2.M (RTE Integrator)
Title : EPA 8270C



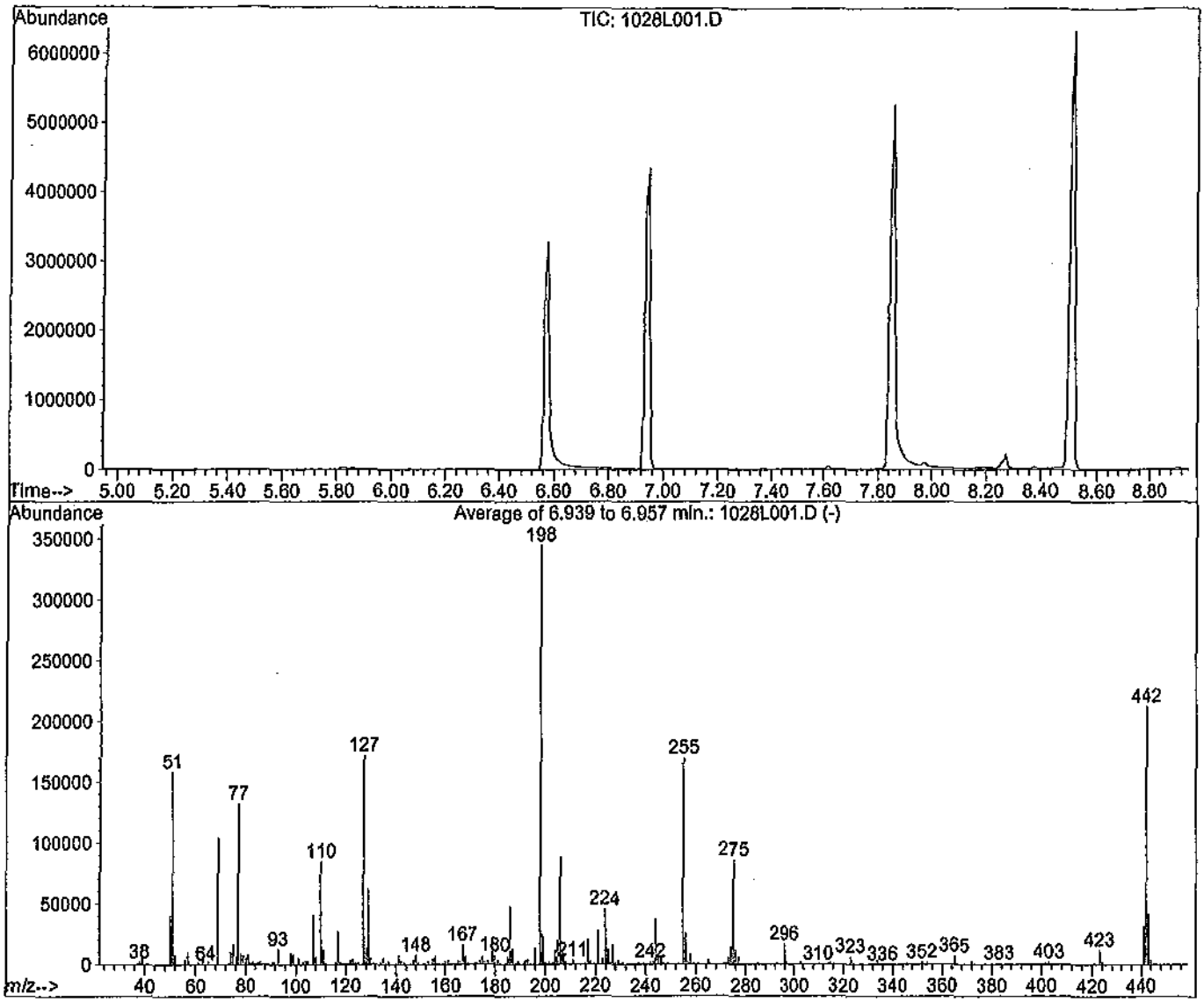
Spectrum Information: Scan 471

Target Mass	Rel. to Mass	Lower Limit%	Upper Limit%	Rel. Abn%	Raw Abn	Result Pass/Fail
51	198	30	60	50.7	294016	PASS
68	69	0.00	2	0.0	0	PASS
70	69	0.00	2	0.6	1188	PASS
127	198	40	60	54.3	314624	PASS
197	198	0.00	1	0.0	0	PASS
198	198	100	100	100.0	579520	PASS
199	198	5	9	7.0	40304	PASS
275	198	10	30	24.5	141888	PASS
365	198	1	100	2.0	11470	PASS
441	443	0.01	100	70.8	44728	PASS
442	198	40	150	52.6	304768	PASS
443	442	17	23	20.7	63176	PASS

DFTPP

Data File : M:\LINUS\DATA\L111027\1028L001.D Vial: 1
Acq On : 28 Oct 11 9:32 Operator: LF
Sample : SVTUNE 10-27-11 Inst : Linus
Misc : Multiplr: 1.00

Method : M:\LINUS\DATA\L111027\SIM2.M (RTE Integrator)
Title : EPA 8270C



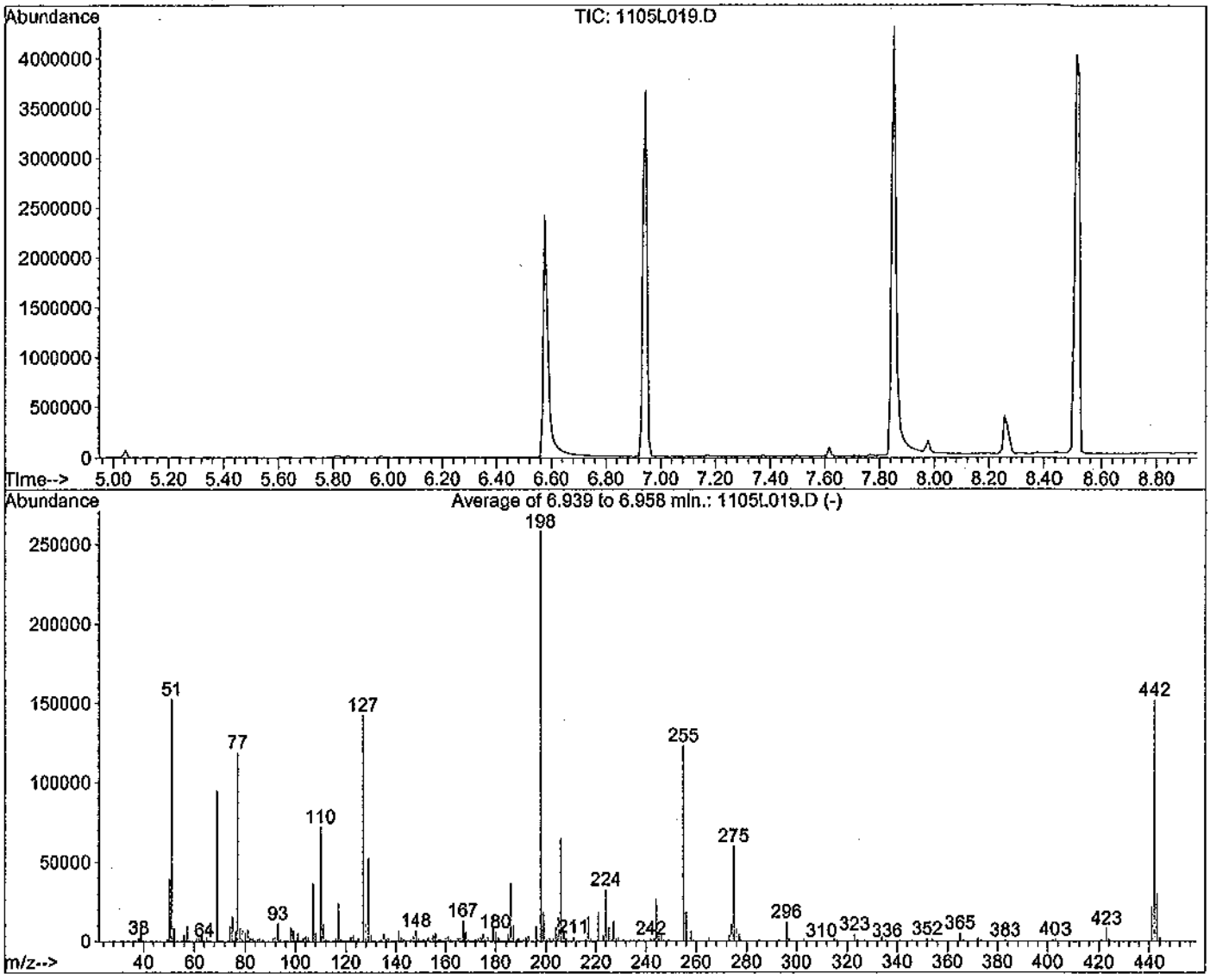
Spectrum Information: Average of 6.939 to 6.957 min.

Target Mass	Rel. to Mass	Lower Limit%	Upper Limit%	Rel. Abn%	Raw Abn	Result Pass/Fail
51	198	30	60	45.8	158326	PASS
68	69	0.00	2	0.0	0	PASS
70	69	0.00	2	0.5	519	PASS
127	198	40	60	49.8	171922	PASS
197	198	0.00	1	0.0	0	PASS
198	198	100	100	100.0	345360	PASS
199	198	5	9	7.1	24580	PASS
275	198	10	30	24.8	85541	PASS
365	198	1	100	2.0	6987	PASS
441	443	0.01	100	74.7	31248	PASS
442	198	40	150	61.5	212309	PASS
443	442	17	23	19.7	41843	PASS

DFTPP

Data File : M:\LINUS\DATA\L111027\1105L019.D Vial: 19
Acq On : 5 Nov 11 16:36 Operator: LF
Sample : SVTUNE 10-27-11 Inst : Linus
Misc : Multiplr: 1.00

Method : M:\LINUS\DATA\L111027\SIM2.M (RTE Integrator)
Title : EPA 8270C



Spectrum Information: Average of 6.939 to 6.958 min.

Target Mass	Rel. to Mass	Lower Limit%	Upper Limit%	Rel. Abn%	Raw Abn	Result Pass/Fail
51	198	30	60	59.0	152381	PASS
68	69	0.00	2	0.0	0	PASS
70	69	0.00	2	0.6	557	PASS
127	198	40	60	55.1	142318	PASS
197	198	0.00	1	0.5	1197	PASS
198	198	100	100	100.0	258253	PASS
199	198	5	9	7.1	18230	PASS
275	198	10	30	23.2	59874	PASS
365	198	1	100	1.9	4901	PASS
441	443	0.01	100	73.0	21870	PASS
442	198	40	150	58.8	151760	PASS
443	442	17	23	19.7	29958	PASS

GC/MS STANDARD PREPARATION BOOK # J PAGE # 84

VF 1/17/11

PREP DATE:	01-17-11					
8270C Stock/Spike Standard						
Exp:	05-29-11					
		Conc.		Date	CODE:	P
Supplier	ID #	µg/mL	Lot #	Code	Exp.Date	µL
Absolute	10001	2000	032009-28092	01/17/11	03-20-12	1000
Absolute	10001	2000	032009-28091	01/17/11	03-20-12	1000
Absolute	10002	2000	073109-27974	01/17/11	07-31-12	1000
Absolute	10002	2000	073109-27973	01/17/11	07-31-12	1000
Absolute	10004	2000	101509-27979	01/17/11	10-15-14	1000
Absolute	10004	2000	101509-27978	01/17/11	10-15-14	1000
Absolute	10005	2000	061209-27984	01/17/11	06-12-14	1000
Absolute	10005	2000	061209-27983	01/17/11	06-12-14	1000
Absolute	10006	2000	120810-27989	01/17/11	12-08-13	1000
Absolute	10006	2000	120810-27988	01/17/11	12-08-13	1000
Absolute	10007	2000	100909-28010	01/17/11	10-09-14	1000
Absolute	10007	2000	100909-28013	01/17/11	10-09-14	1000
Absolute	10018	2000	073109-27994	01/17/11	07-31-14	1000
Absolute	10018	2000	073109-27993	01/17/11	07-31-14	1000
Absolute	70023	1000	080310-28008	01/17/11	08-03-15	1000
Absolute	70023	1000	080310-28009	01/17/11	08-03-15	1000
Absolute	82705	2000	121010-27999	01/17/11	12-10-13	1000
Absolute	82705	2000	121010-27998	01/17/11	12-10-13	1000
Absolute	94552	2000	052908-28004	01/17/11	05-29-11	1000
Absolute	94552	2000	052908-28003	01/17/11	05-29-11	1000
					Final Vol	20000

VF 4/5/11

PREP DATE:	01-25-11																
8270T STANDARD CURVE																	
Exp:	02-24-11						0.1	0.2	1	5	10	20	40	50	60	80	100
		Conc.		Date													
Supplier	ID #	µg/mL	Lot #	Code	Exp.Date				µL	µL	µL	µL	µL	µL	µL	µL	µL
	8270T Stock	200		12/17/10	05-29-11		0	0	0	5	5	10	20	25	30	40	50
	5.0ug/mL			01/25/11			0	0	20	0	0	0	0	0	0	0	0
	1.0ug/mL			01/25/11			10	20	0	0	0	0	0	0	0	0	0
	Surrogate Stock	VAR	160518-27570	11/11/10	11-11-11		0	0	0	5	5	10	20	25	30	40	50
EM Science	Methylene Chloride		47080				90	80	80	190	90	80	60	50	40	20	0
					Final Vol.				100	200	100	100	100	100	100	100	100

VF 1/25/11

PREP DATE:	01-25-11					
8270 Second Source (SS) 50ug/mL						
						50
		Conc.		Date	CODE:	
Supplier	ID #	µg/mL	Lot #	Code	Exp.Date	µL
	8270C SS	200		10/06/10	10-06-11	25
EM Science	Methylene Chloride		47080			75
					Final Vol.	100

VF 1/20/11

Method 8270 Internal Standard Solution, 2,000 mg/L, 1 ml
110001-02

Lot #	Storage	Expiry
167766	5-10 Degrees C	4/20/13

Solvent: Methylene Chloride

8270 Internal Standard
Lot #: 167766 - 28148
Rec: 1/20/11 MFR exp. 04/20/13

agp 1/25/12

VF 1/25/11

Method 8270 Internal Standard Solution, 2,000 mg/L, 1 ml
110001-02

Lot #	Storage	Expiry
167766	5-10 Degrees C	4/20/13

Solvent: Methylene Chloride

8270 Internal Standard
Lot #: 167766 - 28147
Rec: 1/20/11 MFR exp. 04/20/13

agp 1/25/12

GC/MS STANDARD PREPARATION BOOK # 5 PAGE # 89

LF 3/23/11

Part #: 94552 Laboratory Use Only-See MSDS
Lot #: 052908 Exp: 052911 1 mL
Semi-Volatile Standard
11 components
Varied ug/mL in
ABSOLUTE STANDA
Semi-Volatile Standard
Lot #: 052908 - 28001
Rec: 12/16/10 MFR exp. 05/29/11

LF exp 5/29/11

LF 3/23/11

Part #: 94552 Laboratory Use Only-See MSDS
Lot #: 052908 Exp: 052911 1 mL
Semi-Volatile Standard
11 components
Varied ug/mL in
ABSOLUTE STANDA
Semi-Volatile Standard
Lot #: 052908 - 28002
Rec: 12/16/10 MFR exp. 05/29/11

exp 5/29/11

LF 3/23/11

Part #: 82705 Laboratory Use Only - See MSDS
Lot #: 121010 Exp: 121013 Storage 4 °C
EPA Method 8270A
4 components
2000 ug/mL in ace
EPA Method 8270A-Mix#11
Lot #: 121010 - 27996
Rec: 12/16/10 MFR exp. 12/10/13
ABSOLUTE STANDARDS, INC.

LF exp 5/29/11

LF 3/23/11

Part #: 82705 Laboratory Use Only - See MSDS
Lot #: 121010 Exp: 121013 Storage 4 °C
EPA Method 8270A - Mix #11
4 components
2000 ug/mL in ace
EPA Method 8270A-Mix#11
Lot #: 121010 - 27997
Rec: 12/16/10 MFR exp. 12/10/13
ABSOLUTE STANDAF

LF exp 5/29/11

LF 3/23/11

PREP DATE:	03-23-11					
8270C Stock/Spike Standard						
Exp:	05-29-11					
		Conc.		Date	CODE:	P
Supplier	ID #	µg/mL	Lot #	Code	Exp.Date	µL
Absolute	10001	2000	032009-28089	03/23/11	03-20-12	1000
Absolute	10001	2000	320009-28090	03/23/11	03-20-12	1000
Absolute	10002	2000	073109-27971	03/23/11	07-31-12	1000
Absolute	10002	2000	073109-27972	03/23/11	07-31-12	1000
Absolute	10004	2000	101509-27976	03/23/11	10-15-14	1000
Absolute	10004	2000	101509-27977	03/23/11	10-15-14	1000
Absolute	10005	2000	061209-27981	03/23/11	06-12-14	1000
Absolute	10005	2000	061209-27982	03/23/11	06-12-14	1000
Absolute	10006	2000	120810-27986	03/23/11	12-08-13	1000
Absolute	10006	2000	120810-27987	03/23/11	12-08-13	1000
Absolute	10007	2000	100909-28015	03/23/11	10-09-14	1000
Absolute	10007	2000	100909-28014	03/23/11	10-09-14	1000
Absolute	10018	2000	073109-27991	03/23/11	07-31-14	1000
Absolute	10018	2000	073109-27992	03/23/11	07-31-14	1000
Absolute	70023	1000	080310-28006	03/23/11	08-03-15	1000
Absolute	70023	1000	080310-28007	03/23/11	08-03-15	1000
Absolute	82705	2000	052908-28001	03/23/11	05-29-11	1000
Absolute	82705	2000	052908-28002	03/23/11	05-29-11	1000
Absolute	94552	2000	121010-27996	03/23/11	12-10-13	1000
Absolute	94552	2000	121010-27997	03/23/11	12-10-13	1000
					Final Vol	20000

LF 3/23/11

9IM IS exp 1/25/12
1500µL EM Science MC Lot #47088825
100µL 8270 IS opened 1/25/11 exp 1/25/12

GC/MS STANDARD PREPARATION BOOK # J PAGE # 90

VF 3/28/11

o2si smart solutions — 8270 BN:A (200:400) Surrogate Solution, 1 ml
110004-17 — Storage: <=-10 Degrees C
Made in USA — Lot No: 160538 — Solvent: Methylene Chloride
Exp: 4/10/2013
Date Opened:
8270 BN:A (200:400) Surrogate Solution
Lot #: 160538 - 27574
Rec: 10/18/10 MFR exp. 08/10/12

VF exp 3/28/12

VF 3/28/11

PREP DATE:	03-28-11															
8270T STANDARD CURVE																
Exp:	04-27-11						0.1	0.2	1	5	10	20	40	50	60	80	100
		Conc.		Date													
Supplier	ID #	µg/mL	Lot #	Code	Exp.Date				µL	µL	µL	µL	µL	µL	µL	µL	µL
	8270T Stock	200		03/23/11	05-29-11	0	0	0	5	5	10	20	25	30	40	50
	5.0ug/mL			03/28/11		0	0	20	0	0	0	0	0	0	0	0
	1.0ug/mL			03/28/11		10	20	0	0	0	0	0	0	0	0	0
	Surrogate Stock	VAR	160538-27574	03/28/11	03-28-12	0	0	0	5	5	10	20	25	30	40	50
EM Science	Methylene Chloride		47080			90	80	80	190	90	80	60	50	40	20	0
					Final Vol.			100	200	100	100	100	100	100	100	100

VF 3/28/11

PREP DATE:	03-28-11					
8270 Second Source (SS) 50ug/mL						50 VF
		Conc.		Date	CODE:	
Supplier	ID #	µg/mL	Lot #	Code	Exp.Date	µL
	8270C SS	200		10/06/10	10-06-11	25
EM Science	Methylene Chloride		47080			75
					Final Vol.	100

VF 4/18/11

GCM-150-1
Lot: CF-2995
Exp: 08/31/2011
Semi-Volatiles GC/MS Tuning Standard
4 analyte(s) at 1000 µg/mL in dichloromethane
250 Smith St, No Kingstown, RI 02852 USA
ULTRA 1 mL
Semi-volatiles GC/MS Tuning Standard
Lot #: CF-2995 - 26131
Rec: 2/17/10 MFR exp. 08/31/11

VF exp 8/31/11

VF 4/13/11

PREP DATE:	04-13-11						
SV Tune Mix 50ug/ml							
Exp:	08-31-11						
		Conc.			Date	CODE:	B
Supplier	ID #	µg/mL	Lot #		Code	Exp.Date	µL
U. Scientific	GCM-150	1000	CF-2995-26131		04/13/11	08-31-11	1000
EM Science	MeCl2		47080				19000
						Final Vol	20000

exp 8/31/11

VF 4/20/11

8270D PAH SIM Solution, 200 mg/L, 1 ml
110780-01
Lot # 170253 — Storage <-10 Degrees C — Expiry 3/3/13
Solv Methylene Chloride
Made in the USA

VF exp 4/20/12

8270D PAH SIM
Lot #: 170253 - 28485
Rec: 3/10/11 MFR exp. 3/3/2013

VF 4/19/11

8270D PAH SIM Solution, Second Source, 200 mg/L, 1 ml
110780-01-SS
Lot # 170256 — Storage <-10 Degrees C — Expiry 3/3/13
Solv Methylene Chloride
Made in the USA

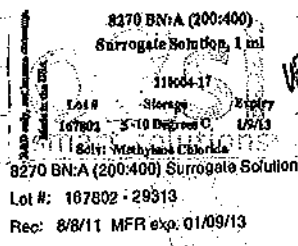
VF exp 4/20/12

8270D PAH SIM (SS)
Lot #: 170256 - 28487
Rec: 3/10/11 MFR exp. 3/3/2013

VF 8/16/11

PREP DATE:	08/16/11		exp:	08/23/11			
10ug/mL 1,2,3-TCP							
	50uL of 1000ug/mL 1,2,3 TCP into a final volume of 5mL of P&T Methanol						
	1000ug/mL 1,2,3 TCP date code:			05/27/11			
	P & T Methonal Lot #			9077-02			
PREP DATE:	08/16/11		exp:	08/23/11			
1ug/mL 1,2,3-TCP							
	5uL of 1000ug/mL 1,2,3 TCP into a final volume of 5mL of P&T Methanol						
	1000ug/mL 1,2,3 TCP date code:			05/27/11			
	P & T Methanol Lot #			JT Baker H46E44			
PREP DATE:	08/16/11		exp:	08/23/11			
2ug/mL 1,2,3-TCPd5							
	10uL of 2000ug/mL 1,2,3 TCP into a final volume of 10mL of P&T Methanol						
	2000ug/mL 1,2,3 TCP-d5 date code:			05/27/11			
	P & T Methanol Lot #			9077-02			

VF 8/22/11



VF exp 8/22/12

VF 8/22/11

PREP DATE:	08-22-11												
8270 STANDARD CURVE													
Exp:	08-29-11						5	10	20	40	50	60	80	100
Supplier	ID #	Conc. µg/mL	Lot #	Date Code	Exp.Date		µL	µL	µL	µL	µL	µL	µL	µL
	8270T Stock	200		07/26/11	01-26-12		5	5	10	20	25	30	40	50
	Surrogate Stock	VAR	167802-29313	08/22/11	08-22-12		5	5	10	20	25	30	40	50
EM Science	Methylene Chloride		47186				190	90	80	60	50	40	20	0
						Final Vol.	200	100	100	100	100	100	100	100

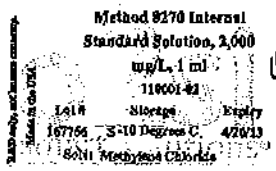
VF 8/22/11

PREP DATE:	08-22-11					
8270 Second Source (SS) 50ug/mL						
						50
Supplier	ID #	Conc. µg/mL	Lot #	Date Code	CODE: Exp.Date	µL
	8270C SS	200		10/06/10	10-06-11	25
EM Science	Methylene Chloride		47186			75
					Final Vol.	100

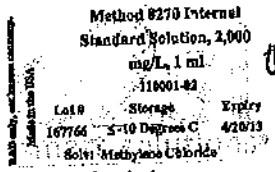
VF 8/24/11

PREP DATE:	09-21-11													
8270 SIM STANDARD CURVE														
							0.10	0.20	0.50	1.00	5.00	10.00	50.00	100.00
Supplier	ID #	Conc. µg/mL	Lot #	Date Code	CODE: Exp.Date		A µL	A µL	C µL	D µL	E µL	F µL	G µL	H µL
	8270D PAH SIM	200	170253-28485	04/20/11	04-20-12		0	0	0	0	5	5	25	50
	5.0ug/mL	5		09/21/11			0	0	10	20	0	0	0	0
	1.0ug/mL	1		09/21/11			10	20	0	0	0	0	0	0
	Surrogate Stock	VAR	167802-29313	08/22/11	08-23-11		0	0	0	0	5	5	25	50
EM Science	Methylene Chloride		47186				90	80	90	80	190	90	50	0
						Final Vol.	100	100	100	100	200	100	100	100

LF 10/18/11



8270 Internal Standard
Lot #: 167766 - 28149
Rec: 1/20/11 MFR exp. 04/20/13



8270 Internal Standard
Lot #: 167766 - 28150
Rec: 1/20/11 MFR exp. 04/20/13

exp 10/18/12

LF 10/27/11

GCM-160-1
Lot: CH-2137
Exp: 07/31/2013
ULTRA 1 mL
Semi-Volatiles GC/MS Tuning Standard
4 analyte(s) at 1000 µg/mL in dichloromethane
250 Smith St, No Kingstown, RI 02852 USA
For Lab Use Only

exp 10/27/12

50µg/mL SV Tune Mix: 1mL of GCM-160-1 Lot# CH2137 into 19mL of Eur Science MC Lot# 42080.

LF 10/27/11

PREP DATE:	10-27-11												
8270 SIM STANDARD CURVE													
						0.10	0.20	0.50	1.00	5.00	10.00	50.00	100.00
		Conc.		Date	CODE:	A	A	C	D	E	F	G	H
Supplier	ID #	µg/mL	Lot #	Code	Exp.Date	µL	µL	µL	µL	µL	µL	µL	µL
	8270D PAH SIM	200	170253-28485	04/20/11	04-20-12	0	0	0	0	5	5	25	50
	5.0ug/mL	5		10/27/11		0	0	10	20	0	0	0	0
	1.0ug/mL	1		10/27/11		10	20	0	0	0	0	0	0
	Surrogate Stock	VAR	167802-29313	08/22/11	08-22-12	0	0	0	0	5	5	25	50
EM Science	Methylene Chloride		47186			90	80	90	80	190	90	50	0
					Final Vol.	100	100	100	100	200	100	100	100

LF 10/27/11

PREP DATE:	10-27-11					
SIM 8270 Second Source (5µg/mL)						
Exp:	11-10-11					
				Conc.	Date	CODE:	
Supplier	ID #		Lot #	µg/mL	Code	Exp.Date	µL
	8270D PAH SIM (SS)		170256-28487	200	04/20/11	04-20-12	5
	MeCl2			Lot#47186			195
					Final Volume		200

LF 11/8/11

PREP DATE:	11-08-11													
8270 STANDARD CURVE														
Exp:	11-15-11						5	10	20	40	50	60	80	100
		Conc.		Date										
Supplier	ID #	µg/mL	Lot #	Code	Exp.Date		µL	µL	µL	µL	µL	µL	µL	µL
	8270T Stock	200		10/18/11	04-18-12		5	5	10	20	25	30	40	50
	Surrogate Stock	VAR	167802-29313	08/22/11	08-22-12		5	5	10	20	25	30	40	50
EM Science	Methylene Chloride		47186				190	90	80	60	50	40	20	0
					Final Vol.		200	100	100	100	100	100	100	100

LF 11/8/11

PREP DATE:	11-08-11					
8270 Second Source (SS) 50ug/mL						
						50
		Conc.		Date	CODE:	
Supplier	ID #	µg/mL	Lot #	Code	Exp.Date	µL
	8270C SS	200		10/11/11	04-12-12	25
EM Science	Methylene Chloride		47186			75
					Final Vol.	100

Organic Extraction Worksheet

Method	SIM Separatory Funnel Extra 3510C	Extraction Set	111031A	Extraction Method	SEP004S	Units	mL

Spiked ID 1	SIM Spike 178987-29587		Surrogate ID 1	8270 SIM Surrogate 172835-28827	
Spiked ID 2			Surrogate ID 2		
Spiked ID 3			Surrogate ID 3		
Spiked ID 4			Surrogate ID 4		
Spiked ID 5			Surrogate ID 5		
Spiked ID 6			Sufficient Vol for Matrix QC:	YES	
Spiked ID 7			Ext. Start Time:		
Spiked ID 8			Ext. End Time:		
			GC Requires Extract By:	11/02/11 0:00	

pH1	2	31/2011 11:25:00 AM	Water Bath Temp Criteria	80 °C
pH2	14	0/31/2011 4:00:00 PM		
pH3				

Spiked By: HW **Date** 10/31/2011 **Witnessed By:** DL **Date** 10/31/2011

	Sample	Sample Container	Spike Amount	Spike ID	Surrogate Amount	Surrogate ID	Extract Amount	Final Volume	pH	Extract Date/Time	Comments
1	111031A Blk				0.025	1	1000	1	2/1	10/31/11 11:20	
						equip	E-WB5				
2	111031A LCS-1		0.025	1	0.025	1	1000	1	2/1	10/31/11 11:20	
						equip	E-WB5				
3	AY49327	AY49327W08			0.025	1	1040	1	2/1	10/31/11 11:20	66103-1 WEEK RUSH -- Amber Liter
						equip	E-WB5				
4	AY49328	AY49328W04			0.025	1	1040	1	2/1	10/31/11 11:20	66103-1 WEEK RUSH -- Amber Liter
						equip	E-WB5				
5	AY49329	AY49329W04			0.025	1	1030	1	2/1	10/31/11 11:20	66103-1 WEEK RUSH -- Amber Liter
						equip	E-WB5				
6	AY49330	AY49330W07			0.025	1	1040	1	2/1	10/31/11 11:20	66103-1 WEEK RUSH -- Amber Liter
						equip	E-WB5				
7	AY49331	AY49331W07			0.025	1	1030	1	2/1	10/31/11 11:20	66103-1 WEEK RUSH -- Amber Liter
						equip	E-WB5				
8	AY49333	AY49333W10			0.025	1	1050	1	2/1	10/31/11 11:20	66102-2 WEEK RUSH -- Amber Liter
						equip	E-WB5				
9	AY49334 MS-1	AY49334W30	0.025	1	0.025	1	1030	1	2/1	10/31/11 11:20	66102-2 WEEK RUSH -- Amber Liter
						equip	E-WB5				
10	AY49334 MSD-1	AY49334W34	0.025	1	0.025	1	1030	1	2/1	10/31/11 11:20	66102-2 WEEK RUSH -- Amber Liter
						equip	E-WB5				
11	AY49334	AY49334W29			0.025	1	1050	1	2/1	10/31/11 11:20	66102-2 WEEK RUSH -- Amber Liter
						equip	E-WB6				
12	AY49336	AY49336W10			0.025	1	1030	1	2/1	10/31/11 11:20	66102-2 WEEK RUSH -- Amber Liter
						equip	E-WB6				
13	AY49481	AY49481W08			0.025	1	1050	1	2/1	10/31/11 11:20	66116-2 WEEK RUSH -- Amber Liter
						equip	E-WB6				

Solvent and Lot#	
MC	EMD 51204
Na2SO4	3581C501
10N NaOH	10/31/11
1+1 Acid	09/15/11
A. Na2SO4	10/31/11

Extraction COC Transfer	
Extraction lab employee Initials	HW
GC analyst's initials	
Date	11/5/11
Time	802
Refrigerator	Hwang

	Technician's Initials
Scanned By	HW
Sample Preparation	CC
Extraction	HW/DL/JL
Concentration	JL

Modified	10/31/2011 10:48:43 AM

Reviewed By: HW 229 **Date** 11/1/2011

_SEP004S

Organic Extraction Worksheet

Method	SIM Separatory Funnel Extra 3510C	Extraction Set	111031A	Extraction Method	SEP004S	Units	mL

Spiked ID 1	SIM Spike 178987-29587		Surrogate ID 1	8270 SIM Surrogate 172835-28827	
Spiked ID 2			Surrogate ID 2		
Spiked ID 3			Surrogate ID 3		
Spiked ID 4			Surrogate ID 4		
Spiked ID 5			Surrogate ID 5		
Spiked ID 6			Sufficient Vol for Matrix QC:	YES	
Spiked ID 7			Ext. Start Time:		
Spiked ID 8			Ext. End Time:		
			GC Requires Extract By:	11/02/11 0:00	

pH1	2	31/2011 11:25:00 AM	Water Bath Temp Criteria	80 °C
pH2	14	0/31/2011 4:00:00 PM		
pH3				

Spiked By: HW **Date** 10/31/2011 **Witnessed By:** DL **Date** 10/31/2011

	Sample	Sample Container	Spike Amount	Spike ID	Surrogate Amount	Surrogate ID	Extract Amount	Final Volume	pH	Extract Date/Time	Comments
14	AY49482	AY49482W08			0.025	1	1030	1	2/1	10/31/11 11:20	66116-2 WEEK RUSH -- Amber Liter
						equip	E-WB6				

HW 11/1/11

Solvent and Lot#	
MC	EMD 51204
Na2SO4	3581C501
10N NaOH	10/31/11
1+1 Acid	09/15/11
A. Na2SO4	10/31/11

Extraction COC Transfer	
Extraction lab employee Initials	HW
GC analyst's initials	
Date	11/1/11
Time	
Refrigerator	

	Technician's Initials
Scanned By	HW
Sample Preparation	CC
Extraction	HW/DL/JL
Concentration	JL

Modified	10/31/2011 10:48:43 AM

Reviewed By: HW 230 **Date** 11/1/2011

Injection Log

Directory: M:\LINUS\DATA\L111027\

Line	Vial	FileName	Multiplier	SampleName	Misc Info	Injected
1	1	1027L001.D	1	SVTUNE 10-27-11		27 Oct 11 18:29
2	3	1027L003.D	1	0.1ug/ml PAH 10-27-11		27 Oct 11 19:12
3	4	1027L004.D	1	0.2ug/ml PAH		27 Oct 11 19:38
4	1	1028L001.D	1	SVTUNE 10-27-11		28 Oct 11 9:32
5	5	1028L005.D	1	0.5ug/ml PAH		28 Oct 11 11:07
6	6	1028L006.D	1	1.0ug/ml PAH		28 Oct 11 11:32
7	7	1028L007.D	1	5.0ug/ml PAH		28 Oct 11 11:58
8	8	1028L008.D	1	10ug/ml PAH		28 Oct 11 12:23
9	9	1028L009.D	1	50ug/ml PAH		28 Oct 11 12:49
10	10	1028L010.D	1	100ug/ml PAH		28 Oct 11 13:14
11	11	1028L011.D	1	5.0ug/ml SS PAH 10-27-11		28 Oct 11 13:40
12	19	1105L019.D	1	SVTUNE 10-27-11		5 Nov 11 16:36
13	20	1105L020.D	1	5.0ug/ml PAH 10-27-11		5 Nov 11 16:54
14	28	1105L028.D	1	111031A BLK 1/1000		5 Nov 11 20:15
15	29	1105L029.D	1	111031A LCS-1 1/1000		5 Nov 11 20:41
16	35	1105L035.D	0.95238	AY49333W10 1/1050		5 Nov 11 23:11
17	36	1105L036.D	0.97087	AY49334W30 MS-1 1/1030		5 Nov 11 23:36
18	37	1105L037.D	0.97087	AY49334W34 MSD-1 1/1030		6 Nov 11 00:01
19	38	1105L038.D	0.95238	AY49334W29 1/1050		6 Nov 11 00:26
20	39	1105L039.D	0.97087	AY49336W10 1/1030		6 Nov 11 00:51

EPA METHOD 8260B
Volatile Organic Compounds

APPL, INC.

EPA METHOD 8260B
Volatile Organic Compounds
QC Summary

APPL, INC.

Method Blank

EPA 8260B VOCs + Gas Water

Blank Name/QCG: **111030W-49334 - 161029**
Batch ID: #86RHB-111030AC

APPL Inc.
908 North Temperance Avenu
Clovis, CA 93611

Sample Type	Analyte	Result	LOQ	LOD	DL	Units	Extraction Date	Analysis Date
BLANK	1,1,1,2-TETRACHLOROETHANE	0.26 U	1.0	0.26	0.13	ug/L	10/31/11	10/31/11
BLANK	1,1,1-TRICHLOROETHANE	0.28 U	1.0	0.28	0.14	ug/L	10/31/11	10/31/11
BLANK	1,1,2,2-TETRACHLOROETHANE	0.20 U	1.0	0.20	0.10	ug/L	10/31/11	10/31/11
BLANK	1,1,2-TRICHLOROETHANE	0.40 U	1.0	0.40	0.20	ug/L	10/31/11	10/31/11
BLANK	1,1-DICHLOROETHANE	0.38 U	1.0	0.38	0.19	ug/L	10/31/11	10/31/11
BLANK	1,1-DICHLOROETHENE	0.60 U	1.0	0.60	0.30	ug/L	10/31/11	10/31/11
BLANK	1,2,3-TRICHLOROPROPANE	0.78 U	2.0	0.78	0.39	ug/L	10/31/11	10/31/11
BLANK	1,2,4-TRICHLOROBENZENE	0.42 U	1.0	0.42	0.21	ug/L	10/31/11	10/31/11
BLANK	1,2-DIBROMO-3-CHLOROPROPA	1.52 U	2.0	1.52	0.76	ug/L	10/31/11	10/31/11
BLANK	1,2-DIBROMOETHANE	0.40 U	1.0	0.40	0.20	ug/L	10/31/11	10/31/11
BLANK	1,2-DICHLOROBENZENE	0.34 U	1.0	0.34	0.17	ug/L	10/31/11	10/31/11
BLANK	1,2-DICHLOROETHANE	0.28 U	1.0	0.28	0.14	ug/L	10/31/11	10/31/11
BLANK	1,2-DICHLOROPROPANE	0.34 U	1.0	0.34	0.17	ug/L	10/31/11	10/31/11
BLANK	1,3-DICHLOROBENZENE	0.22 U	1.0	0.22	0.11	ug/L	10/31/11	10/31/11
BLANK	1,3-DICHLOROPROPENE, TOTAL	0.36 U	1.0	0.36	0.18	ug/L	10/31/11	10/31/11
BLANK	1,4-DICHLOROBENZENE	0.38 U	1.0	0.38	0.19	ug/L	10/31/11	10/31/11
BLANK	2-BUTANONE	1.20 U	10.0	1.20	0.60	ug/L	10/31/11	10/31/11
BLANK	4-METHYL-2-PENTANONE	3.80 U	10.0	3.80	1.90	ug/L	10/31/11	10/31/11
BLANK	ACETONE	1.90 U	10.0	1.90	0.95	ug/L	10/31/11	10/31/11
BLANK	BENZENE	0.32 U	1.0	0.32	0.16	ug/L	10/31/11	10/31/11
BLANK	BROMODICHLOROMETHANE	0.28 U	1.0	0.28	0.14	ug/L	10/31/11	10/31/11
BLANK	BROMOFORM	0.28 U	1.0	0.28	0.14	ug/L	10/31/11	10/31/11
BLANK	BROMOMETHANE	0.48 U	2.0	0.48	0.24	ug/L	10/31/11	10/31/11
BLANK	CARBON TETRACHLORIDE	0.20 U	1.0	0.20	0.10	ug/L	10/31/11	10/31/11
BLANK	CHLOROBENZENE	0.42 U	1.0	0.42	0.21	ug/L	10/31/11	10/31/11
BLANK	CHLORODIBROMOMETHANE	0.38 U	1.0	0.38	0.19	ug/L	10/31/11	10/31/11
BLANK	CHLOROETHANE	0.42 U	1.0	0.42	0.21	ug/L	10/31/11	10/31/11
BLANK	CHLOROFORM	0.14 U	1.0	0.14	0.07	ug/L	10/31/11	10/31/11
BLANK	CHLOROMETHANE	0.84 U	1.0	0.84	0.42	ug/L	10/31/11	10/31/11
BLANK	CIS-1,2-DICHLOROETHENE	0.32 U	1.0	0.32	0.16	ug/L	10/31/11	10/31/11
BLANK	ETHYLBENZENE	0.46 U	1.0	0.46	0.23	ug/L	10/31/11	10/31/11
BLANK	GASOLINE	12.12 U	20.0	12.12	6.06	ug/L	10/31/11	10/31/11
BLANK	HEXACHLOROBUTADIENE	0.38 U	1.0	0.38	0.19	ug/L	10/31/11	10/31/11
BLANK	METHYL TERT-BUTYL ETHER	0.52 U	1.0	0.52	0.26	ug/L	10/31/11	10/31/11

Quant Method: CALLW.M
Run #: 1030C34
Instrument: Chico
Sequence: C111030
Initials: ARS

GC SC-Blank-REG MDLs
Printed: 12/06/11 3:06:05 PM

Method Blank

EPA 8260B VOCs + Gas Water

Blank Name/QCG: **111030W-49334 - 161029**
Batch ID: #86RHB-111030AC

APPL Inc.
908 North Temperance Avenu
Clovis, CA 93611

Sample Type	Analyte	Result	LOQ	LOD	DL	Units	Extraction Date	Analysis Date
BLANK	METHYLENE CHLORIDE	0.70 U	5.0	0.70	0.35	ug/L	10/31/11	10/31/11
BLANK	STYRENE	0.50 U	1.0	0.50	0.25	ug/L	10/31/11	10/31/11
BLANK	TETRACHLOROETHENE	0.48 U	1.0	0.48	0.24	ug/L	10/31/11	10/31/11
BLANK	TOLUENE	0.34 U	1.0	0.34	0.17	ug/L	10/31/11	10/31/11
BLANK	TRANS-1,2-DICHLOROETHENE	0.38 U	1.0	0.38	0.19	ug/L	10/31/11	10/31/11
BLANK	TRICHLOROETHENE	0.32 U	1.0	0.32	0.16	ug/L	10/31/11	10/31/11
BLANK	VINYL CHLORIDE	0.46 U	1.0	0.46	0.23	ug/L	10/31/11	10/31/11
BLANK	XYLENES (TOTAL)	0.38 U	1.0	0.38	0.19	ug/L	10/31/11	10/31/11
BLANK	SURROGATE: 1,2-DICHLOROET	109	70-120			%	10/31/11	10/31/11
BLANK	SURROGATE: 4-BROMOFLUORO	102	75-120			%	10/31/11	10/31/11
BLANK	SURROGATE: DIBROMOFLUOR	105	85-115			%	10/31/11	10/31/11
BLANK	SURROGATE: TOLUENE-D8 (S)	102	85-120			%	10/31/11	10/31/11

Quant Method: CALLW.M
Run #: 1030C34
Instrument: Chlco
Sequence: C111030
Initials: ARS

GC SC-Blank-REG MDLs
Printed: 12/06/11 3:08:05 PM

Form 2 & 8

Surrogate Recovery

Lab Name: APPL, Inc. SDG No: 66102

Case No: 66102 Date Analyzed: 10/31/11

Matrix: WATER Instrument: Chico

APPL ID.	Client Sample No.	SURROGATE: 1,2-DICHLOROETHANE-D4 (S)			SURROGATE: 4-BROMOFLUOROBENZENE (S)		
		Limits	Result	Qualifier	Limits	Result	Qualifier
111030AC-LCS	Lab Control Spike	70-120	101		75-120	109	
111030AC-BLK	Blank	70-120	109		75-120	102	
AY49335	ES048	70-120	108		75-120	102	
AY49333	ES046	70-120	114		75-120	103	
AY49334	ES047	70-120	104		75-120	98.9	
AY49336	ES049	70-120	98.1		75-120	102	
AY49334-MS	Matrix Spike	70-120	93.7		75-120	101	
AY49334-MSD	Matrix SpikeD	70-120	92.1		75-120	105	

Comments: Batch: #86RHB-111030AC

Form 2 & 8

Surrogate Recovery

Lab Name: APPL, Inc. SDG No: 66102

Case No: 66102 Date Analyzed: 10/31/11

Matrix: WATER Instrument: Chico

APPL ID.	Client Sample No.	SURROGATE: DIBROMOFLUOROMETHANE (S)			SURROGATE: TOLUENE-D8 (S)		
		Limits	Result	Qualifier	Limits	Result	Qualifier
111030AC-LCS	Lab Control Spike	85-115	100		85-120	104	
111030AC-BLK	Blank	85-115	105		85-120	102	
AY49335	ES048	85-115	105		85-120	99.0	
AY49333	ES046	85-115	105		85-120	98.9	
AY49334	ES047	85-115	102		85-120	97.6	
AY49336	ES049	85-115	100		85-120	101	
AY49334-MS	Matrix Spike	85-115	98.4		85-120	102	
AY49334-MSD	Matrix SpikeD	85-115	99.6		85-120	104	

Comments: Batch: #86RHB-111030AC

Laboratory Control Spike Recovery
EPA 8260B VOCs + Gas Water

APPL ID: **111031W-49334 LCS - 161029**
Batch ID: #86RHB-111030AC

APPL Inc.
908 North Temperance Avenue
Clovis, CA 93611

Compound Name	Spike Level ug/L	SPK Result ug/L	SPK % Recovery	Recovery Limits
1,1,1,2-TETRACHLOROETHANE	10.00	11.0	110	80-130
1,1,1-TRICHLOROETHANE	10.00	9.55	95.5	65-130
1,1,2,2-TETRACHLOROETHANE	10.00	10.3	103	65-130
1,1,2-TRICHLOROETHANE	10.00	9.91	99.1	75-125
1,1-DICHLOROETHANE	10.00	9.98	99.8	70-135
1,1-DICHLOROETHENE	10.00	8.83	88.3	70-130
1,2,3-TRICHLOROPROPANE	10.00	11.8	118	75-125
1,2,4-TRICHLOROBENZENE	10.00	10.2	102	65-135
1,2-DIBROMO-3-CHLOROPROPANE	10.00	10.2	102	50-130
1,2-DIBROMOETHANE	10.00	10.7	107	70-130
1,2-DICHLOROBENZENE	10.00	10.1	101	70-120
1,2-DICHLOROETHANE	10.00	9.61	96.1	70-130
1,2-DICHLOROPROPANE	10.00	10.1	101	75-125
1,3-DICHLOROBENZENE	10.00	9.67	96.7	75-125
1,3-DICHLOROPROPENE, TOTAL	20.0	19.7	98.5	70-130
1,4-DICHLOROBENZENE	10.00	9.94	99.4	75-125
2-BUTANONE	10.00	10.5	105	30-150
4-METHYL-2-PENTANONE	10.00	9.78	97.8	60-135
ACETONE	10.00	13.0	130	40-140
BENZENE	10.00	9.48	94.8	80-120
BROMODICHLOROMETHANE	10.00	10.0	100	75-120
BROMOFORM	10.00	9.32	93.2	70-130
BROMOMETHANE	10.00	8.95	89.5	30-145
CARBON TETRACHLORIDE	10.00	9.77	97.7	65-140
CHLOROBENZENE	10.00	10.3	103	80-120
CHLORODIBROMOMETHANE	10.00	11.0	110	60-135

Comments:

Primary	SPK
Quant Method :	CALLW.M
Extraction Date :	10/31/11
Analysis Date :	10/31/11
Instrument :	Chico
Run :	1030C28
Initials :	ARS

Printed: 12/06/11 3:06:11 PM

APPL Standard LCS

Laboratory Control Spike Recovery
EPA 8260B VOCs + Gas Water

APPL ID: **111031W-49334 LCS - 161029**
Batch ID: #86RHB-111030AC

APPL Inc.
908 North Temperance Avenue
Clovis, CA 93611

Compound Name	Spike Level ug/L	SPK Result ug/L	SPK % Recovery	Recovery Limits
CHLOROETHANE	10.00	8.89	88.9	60-135
CHLOROFORM	10.00	9.79	97.9	65-135
CHLOROMETHANE	10.00	9.42	94.2	40-125
CIS-1,2-DICHLOROETHENE	10.00	9.18	91.8	70-125
ETHYLBENZENE	10.00	10.1	101	75-125
GASOLINE	300	370	123	75-125
HEXACHLOROBUTADIENE	10.00	10.5	105	50-140
METHYL TERT-BUTYL ETHER	10.00	9.69	96.9	65-125
METHYLENE CHLORIDE	10.00	9.05	90.5	55-140
STYRENE	10.00	10.6	106	65-135
TETRACHLOROETHENE	10.00	10.1	101	45-150
TOLUENE	10.00	9.40	94.0	75-120
TRANS-1,2-DICHLOROETHENE	10.00	9.05	90.5	60-140
TRICHLOROETHENE	10.00	9.64	96.4	70-125
VINYL CHLORIDE	10.00	10.2	102	50-145
XYLENES (TOTAL)	30.0	30.3	101	80-120
SURROGATE: 1,2-DICHLOROETHANE-D	24.2	24.4	101	70-120
SURROGATE: 4-BROMOFLUOROBENZE	25.5	27.7	109	75-120
SURROGATE: DIBROMOFLUOROMETH	25.1	25.2	100	85-115
SURROGATE: TOLUENE-D8 (S)	25.8	26.9	104	85-120

Comments: _____

Primary	SPK
Quant Method :	CALLW.M
Extraction Date :	10/31/11
Analysis Date :	10/31/11
Instrument :	Chico
Run :	1030C28
Initials :	ARS

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APPL Standard LCS

Matrix Spike Recoveries

EPA 8260B VOCs + Gas Water

APPL ID: **111031W-49334 MS - 161029**
Batch ID: #86RHB-111030AC
Sample ID: AY49334
Client ID: ES047

APPL Inc.
908 North Temperance Avenue
Clovis, CA 93611

Compound Name	Spike Lvl ug/L	Matrix Result ug/L	SPK Result ug/L	DUP Result ug/L	SPK % Recovery	DUP % Recovery	Recovery Limits	RPD %	RPD Limits
1,1,1,2-TETRACHLOROETHANE	10.00	ND	10.9	11.2	109	112	80-130	2.7	30
1,1,1-TRICHLOROETHANE	10.00	ND	8.95	9.51	89.5	95.1	65-130	6.1	30
1,1,2,2-TETRACHLOROETHANE	10.00	ND	12.7	12.6	127	126	65-130	0.79	30
1,1,2-TRICHLOROETHANE	10.00	ND	10.6	10.6	106	106	75-125	0.0	30
1,1-DICHLOROETHANE	10.00	ND	9.46	9.94	94.6	99.4	70-135	4.9	30
1,1-DICHLOROETHENE	10.00	ND	8.96	9.50	89.6	95.0	70-130	5.9	30
1,2,3-TRICHLOROPROPANE	10.00	ND	11.6	11.4	116	114	75-125	1.7	30
1,2,4-TRICHLOROBENZENE	10.00	ND	9.55	10.1	95.5	101	65-135	5.6	30
1,2-DIBROMO-3-CHLOROPROPANE	10.00	ND	11.2	12.2	112	122	50-130	8.5	30
1,2-DIBROMOETHANE	10.00	ND	10.9	10.8	109	108	70-130	0.92	30
1,2-DICHLOROBENZENE	10.00	ND	9.85	10.2	98.5	102	70-120	3.5	30
1,2-DICHLOROETHANE	10.00	ND	8.97	9.48	89.7	94.8	70-130	5.5	30
1,2-DICHLOROPROPANE	10.00	ND	9.62	10.5	96.2	105	75-125	8.7	30
1,3-DICHLOROBENZENE	10.00	ND	9.68	10.1	96.8	101	75-125	4.2	30
1,3-DICHLOROPROPENE, TOTAL	20.0	ND	20.4	20.9	102	105	70-130	2.4	30
1,4-DICHLOROBENZENE	10.00	ND	9.80	10.2	98.0	102	75-125	4.0	30
2-BUTANONE	10.00	ND	10.5	10.2	105	102	30-150	2.9	30
4-METHYL-2-PENTANONE	10.00	ND	11.0	11.0	110	110	60-135	0.0	30
ACETONE	10.00	ND	11.9	13.6	119	136	40-140	13.3	30
BENZENE	10.00	ND	9.56	10.2	95.6	102	80-120	6.5	30
BROMODICHLOROMETHANE	10.00	ND	9.85	10.3	98.5	103	75-120	4.5	30
BROMOFORM	10.00	ND	9.59	10.1	95.9	101	70-130	5.2	30
BROMOMETHANE	10.00	ND	9.35	9.85	93.5	98.5	30-145	5.2	30
CARBON TETRACHLORIDE	10.00	ND	9.77	9.93	97.7	99.3	65-140	1.6	30
CHLOROBENZENE	10.00	ND	9.68	10.1	96.8	101	80-120	4.2	30

= Recovery is outside QC limits.

Comments:

Primary	SPK	DUP
Quant Method :	CALLW.M	CALLW.M
Extraction Date :	10/31/11	10/31/11
Analysis Date :	10/31/11	10/31/11
Instrument :	Chico	Chico
Run :	1030C39	1030C40
Initials :	ARS	

Printed: 12/06/11 3:06:12 PM
APPL MSD SCII

Matrix Spike Recoveries

EPA 8260B VOCs + Gas Water

APPL ID: **111031W-49334 MS - 161029**
Batch ID: #86RHB-111030AC
Sample ID: AY49334
Client ID: ES047

APPL Inc.
908 North Temperance Avenue
Clovis, CA 93611

Compound Name	Spike Lvl ug/L	Matrix Result ug/L	SPK Result ug/L	DUP Result ug/L	SPK % Recovery	DUP % Recovery	Recovery Limits	RPD %	RPD Limits
CHLORODIBROMOMETHANE	10.00	ND	10.8	11.4	108	114	60-135	5.4	30
CHLOROETHANE	10.00	ND	8.88	9.26	88.8	92.6	60-135	4.2	30
CHLOROFORM	10.00	ND	9.21	9.65	92.1	96.5	65-135	4.7	30
CHLOROMETHANE	10.00	ND	8.69	9.03	86.9	90.3	40-125	3.8	30
CIS-1,2-DICHLOROETHENE	10.00	ND	9.19	9.57	91.9	95.7	70-125	4.1	30
ETHYLBENZENE	10.00	ND	9.73	10.3	97.3	103	75-125	5.7	30
GASOLINE	300	ND	374	377	125	126 #	75-125	0.80	30
HEXACHLOROBUTADIENE	10.00	ND	9.04	9.78	90.4	97.8	50-140	7.9	30
METHYL TERT-BUTYL ETHER	10.00	ND	10.1	10.2	101	102	65-125	0.99	30
METHYLENE CHLORIDE	10.00	ND	9.83	9.84	98.3	98.4	55-140	0.10	30
STYRENE	10.00	ND	10.2	10.6	102	106	65-135	3.8	30
TETRACHLOROETHENE	10.00	ND	9.48	10.3	94.8	103	45-150	8.3	30
TOLUENE	10.00	ND	9.46	10.4	94.6	104	75-120	9.5	30
TRANS-1,2-DICHLOROETHENE	10.00	ND	9.17	9.47	91.7	94.7	60-140	3.2	30
TRICHLOROETHENE	10.00	ND	9.00	9.38	90.0	93.8	70-125	4.1	30
VINYL CHLORIDE	10.00	ND	9.90	9.78	99.0	97.8	50-145	1.2	30
XYLENES (TOTAL)	30.0	ND	29.1	30.7	97.0	102	80-120	5.4	30
SURROGATE: 1,2-DICHLOROETHANE-D	24.2	NA	22.7	22.3	93.7	92.1	70-120		
SURROGATE: 4-BROMOFLUOROBENZE	25.5	NA	25.8	26.7	101	105	75-120		
SURROGATE: DIBROMOFLUOROMETH	25.1	NA	24.7	25.0	98.4	99.6	85-115		
SURROGATE: TOLUENE-D8 (S)	25.8	NA	26.3	26.9	102	104	85-120		

= Recovery is outside QC limits.

Comments:

Primary	SPK	DUP
Quant Method :	CALLW.M	CALLW.M
Extraction Date :	10/31/11	10/31/11
Analysis Date :	10/31/11	10/31/11
Instrument :	Chico	Chico
Run :	1030C39	1030C40
Initials :	ARS	

Printed: 12/06/11 3:06:12 PM
APPL MSD SCII

EPA 8260B

Form 4

Blank Summary

Lab Name: APPL, Inc. SDG No: 66102

Case No: 66102 Date Analyzed: 10/31/11

Matrix: WATER Instrument: Chico

Blank ID: 111030AC-BLK Time Analyzed: 1302

APPL ID.	Client Sample No.	File ID.	Date Analyzed
111030AC-LCS	Lab Control Spike	1030C28	10/31/11 0848
111030AC-BLK	Blank	1030C34	10/31/11 1302
AY49335	ES048	1030C35	10/31/11 1339
AY49333	ES046	1030C36	10/31/11 1416
AY49334	ES047	1030C37	10/31/11 1453
AY49336	ES049	1030C38	10/31/11 1531
111030AC-MS	Matrix Spike	1030C39	10/31/11 1608
111030AC-MSD	Matrix SpikeD	1030C40	10/31/11 1645

Comments: Batch: #86RHB-111030AC

Form 5
Tune Summary

Lab Name: APPL Inc.
Case No: 66102
Matrix: Water
ID: 20ug/mL BFB STD 10-19-11

SDG No: 66102
Date Analyzed: 10/31/11
Instrument: Chico
Time Analyzed: 7:21

	Client Sample No.	APPL ID.	File ID.	Date Analyzed
1	Lab Control Spike	111030A LCS-1WC (SS)	1030C28W.D	10/31/11 8:48
2	Lab Control Spike	GAS 300ug/L LCS-1WC	1030C31W.D	10/31/11 10:57
3	Blank	111030A BLk-1WC	1030C34W.D	10/31/11 13:02
4	ES048	AY49335W01	1030C35W.D	10/31/11 13:39
5	ES046	AY49333W04	1030C36W.D	10/31/11 14:16
6	ES047	AY49334W13	1030C37W.D	10/31/11 14:53
7	ES049	AY49336W04	1030C38W.D	10/31/11 15:31
8	Matrix Spike	AY49334W141516 MS-1W	1030C39W.D	10/31/11 16:08
9	Matrix Spike Dup	AY49334W141516 MSD-1	1030C40W.D	10/31/11 16:45
10	Matrix Spike	AY49334W161718 MS-1W	1030C41W.D	10/31/11 17:22
11	Matrix Spike Dup	AY49334W161718 MSD-1	1030C42W.D	10/31/11 17:59
12				
13				
14				
15				
16				
17				
18				
19				
20				
21				
22				

m/e		
50	15 - 40% of mass 95	20.6
75	30 - 60% of mass 95	49.7
95	100 - 100% of mass 95	100.0
96	5 - 9% of mass 95	6.6
173	0 - 2% of mass 174	0.0
174	50 - 100% of mass 95	91.6
175	5 - 9% of mass 174	6.9
176	95 - 101% of mass 174	97.7
177	5 - 9% of mass 176	6.7

8A
INTERNAL STANDARD AREA AND RT SUMMARY

Lab Name: APPL Inc. Contract: Review

Lab Code: SDG No.: 66102

Lab File ID (Standard): 1030C20W.D Date Analyzed: 10/31/11

Instrument ID: Chico Time Analyzed: 3:03

GC Column: ID: Heated Purge: (Y/N)

		Fluorobenzene (IS)		Chlorobenzene-D5 (IS)		1,4-Dichlorobenzene-D (IS)	
		AREA #	RT #	AREA #	RT #	AREA #	RT #
	12 HOUR STD	556544	12.84	375296	18.04	203520	22.24
	UPPER LIMIT	1113088	13.34	750592	18.54	407040	22.74
	LOWER LIMIT	278272	12.34	187648	17.54	101760	21.74
	SAMPLE NO.						
01	111030A LCS-1WC (SS	600576	12.84	389760	18.04	212800	22.24
02	111030A BLk-1WC	564158	12.85	392640	18.05	207424	22.26
03	AY49335W01	576569	12.85	396544	18.05	214592	22.25
04	AY49333W04	558622	12.85	391168	18.05	206336	22.26
05	AY49334W13	617536	12.85	437056	18.05	232960	22.26
06	AY49336W04	662144	12.85	450969	18.05	251840	22.26
07	AY49334W141516 MS-	673315	12.85	444608	18.05	234048	22.26
08	AY49334W141516 MSD	673885	12.85	449984	18.05	240256	22.26
09							
10							
11							
12							
13							
14							
15							
16							
17							
18							
19							
20							
21							
22							

AREA UPPER LIMIT = +100% of internal standard area.
AREA LOWER LIMIT = -50% of internal standard area.
RT UPPER LIMIT = +0.50 minutes of internal standard RT
RT LOWER LIMIT = -0.50 minutes of internal standard RT

Column used to flag values outside QC limits with an asterisk.
* Values outside of QC limits.

EPA METHOD 8260B
Volatile Organic Compounds
Sample Data

APPL, INC.

EPA 8260B VOCs + Gas Water

Environet, Inc.
650 Iwilei Rd, #204
Honolulu, HI 96817

APPL Inc.
908 North Temperance Avenue
Clovis, CA 93611

Attn: Stacy Fineran

Project: RED HILL/1022-024

ARF: 66102

Sample ID: ES046

APPL ID: AY49333

Sample Collection Date: 10/24/11

QCG: #86RHB-111030AC-161029

Method	Analyte	Result	LOQ	LOD	DL	Units	Extraction Date	Analysis Date
EPA 8260B	1,1,1,2-TETRACHLOROETHANE	0.26 U	1.0	0.26	0.13	ug/L	10/31/11	10/31/11
EPA 8260B	1,1,1-TRICHLOROETHANE	0.28 U	1.0	0.28	0.14	ug/L	10/31/11	10/31/11
EPA 8260B	1,1,2,2-TETRACHLOROETHANE	0.20 U	1.0	0.20	0.10	ug/L	10/31/11	10/31/11
EPA 8260B	1,1,2-TRICHLOROETHANE	0.40 U	1.0	0.40	0.20	ug/L	10/31/11	10/31/11
EPA 8260B	1,1-DICHLOROETHANE	0.38 U	1.0	0.38	0.19	ug/L	10/31/11	10/31/11
EPA 8260B	1,1-DICHLOROETHENE	0.60 U	1.0	0.60	0.30	ug/L	10/31/11	10/31/11
EPA 8260B	1,2,3-TRICHLOROPROPANE	0.78 U	2.0	0.78	0.39	ug/L	10/31/11	10/31/11
EPA 8260B	1,2,4-TRICHLOROBENZENE	0.42 U	1.0	0.42	0.21	ug/L	10/31/11	10/31/11
EPA 8260B	1,2-DIBROMO-3-CHLOROPROPANE	1.52 U	2.0	1.52	0.76	ug/L	10/31/11	10/31/11
EPA 8260B	1,2-DIBROMOETHANE	0.40 U	1.0	0.40	0.20	ug/L	10/31/11	10/31/11
EPA 8260B	1,2-DICHLOROBENZENE	0.34 U	1.0	0.34	0.17	ug/L	10/31/11	10/31/11
EPA 8260B	1,2-DICHLOROETHANE	0.28 U	1.0	0.28	0.14	ug/L	10/31/11	10/31/11
EPA 8260B	1,2-DICHLOROPROPANE	0.34 U	1.0	0.34	0.17	ug/L	10/31/11	10/31/11
EPA 8260B	1,3-DICHLOROBENZENE	0.22 U	1.0	0.22	0.11	ug/L	10/31/11	10/31/11
EPA 8260B	1,3-DICHLOROPROPENE, TOTAL	0.36 U	1.0	0.36	0.18	ug/L	10/31/11	10/31/11
EPA 8260B	1,4-DICHLOROBENZENE	0.38 U	1.0	0.38	0.19	ug/L	10/31/11	10/31/11
EPA 8260B	2-BUTANONE	1.20 U	10.0	1.20	0.60	ug/L	10/31/11	10/31/11
EPA 8260B	4-METHYL-2-PENTANONE	3.80 U	10.0	3.80	1.90	ug/L	10/31/11	10/31/11
EPA 8260B	ACETONE	1.90 U	10.0	1.90	0.95	ug/L	10/31/11	10/31/11
EPA 8260B	BENZENE	0.32 U	1.0	0.32	0.16	ug/L	10/31/11	10/31/11
EPA 8260B	BROMODICHLOROMETHANE	0.28 U	1.0	0.28	0.14	ug/L	10/31/11	10/31/11
EPA 8260B	BROMOFORM	0.28 U	1.0	0.28	0.14	ug/L	10/31/11	10/31/11
EPA 8260B	BROMOMETHANE	0.48 U	2.0	0.48	0.24	ug/L	10/31/11	10/31/11
EPA 8260B	CARBON TETRACHLORIDE	0.20 U	1.0	0.20	0.10	ug/L	10/31/11	10/31/11
EPA 8260B	CHLOROBENZENE	0.42 U	1.0	0.42	0.21	ug/L	10/31/11	10/31/11
EPA 8260B	CHLORODIBROMOMETHANE	0.38 U	1.0	0.38	0.19	ug/L	10/31/11	10/31/11
EPA 8260B	CHLOROETHANE	0.42 U	1.0	0.42	0.21	ug/L	10/31/11	10/31/11
EPA 8260B	CHLOROFORM	0.14 U	1.0	0.14	0.07	ug/L	10/31/11	10/31/11
EPA 8260B	CHLOROMETHANE	0.84 U	1.0	0.84	0.42	ug/L	10/31/11	10/31/11
EPA 8260B	CIS-1,2-DICHLOROETHENE	0.32 U	1.0	0.32	0.16	ug/L	10/31/11	10/31/11
EPA 8260B	ETHYLBENZENE	0.46 U	1.0	0.46	0.23	ug/L	10/31/11	10/31/11
EPA 8260B	GASOLINE	12.12 U	20.0	12.12	6.06	ug/L	10/31/11	10/31/11
EPA 8260B	HEXACHLOROBUTADIENE	0.38 U	1.0	0.38	0.19	ug/L	10/31/11	10/31/11
EPA 8260B	METHYL TERT-BUTYL ETHER	0.52 U	1.0	0.52	0.26	ug/L	10/31/11	10/31/11
EPA 8260B	METHYLENE CHLORIDE	0.70 U	5.0	0.70	0.35	ug/L	10/31/11	10/31/11

Quant Method: CALLW.M
Run #: 1030C36
Instrument: Chico
Sequence: C111030
Dilution Factor: 1
Initials: ARS

Printed: 12/06/11 3:06:18 PM

APPL-F1-SC-NoMC-REG MDLs

EPA 8260B VOCs + Gas Water

Environet, Inc.
650 Iwilei Rd, #204
Honolulu, HI 96817

APPL Inc.
908 North Temperance Avenue
Clovis, CA 93611

Attn: Stacy Fineran

Project: RED HILL/1022-024

ARF: 66102

Sample ID: ES046

APPL ID: AY49333

Sample Collection Date: 10/24/11

QCG: #86RHB-111030AC-161029

Method	Analyte	Result	LOQ	LOD	DL	Units	Extraction Date	Analysis Date
EPA 8260B	STYRENE	0.50 U	1.0	0.50	0.25	ug/L	10/31/11	10/31/11
EPA 8260B	TETRACHLOROETHENE	0.48 U	1.0	0.48	0.24	ug/L	10/31/11	10/31/11
EPA 8260B	TOLUENE	0.34 U	1.0	0.34	0.17	ug/L	10/31/11	10/31/11
EPA 8260B	TRANS-1,2-DICHLOROETHENE	0.38 U	1.0	0.38	0.19	ug/L	10/31/11	10/31/11
EPA 8260B	TRICHLOROETHENE	0.32 U	1.0	0.32	0.16	ug/L	10/31/11	10/31/11
EPA 8260B	VINYL CHLORIDE	0.46 U	1.0	0.46	0.23	ug/L	10/31/11	10/31/11
EPA 8260B	XYLENES (TOTAL)	0.38 U	1.0	0.38	0.19	ug/L	10/31/11	10/31/11
EPA 8260B	SURROGATE: 1,2-DICHLOROETHANE-	114	70-120			%	10/31/11	10/31/11
EPA 8260B	SURROGATE: 4-BROMOFLUOROBEN	103	75-120			%	10/31/11	10/31/11
EPA 8260B	SURROGATE: DIBROMOFLUOROMET	105	85-115			%	10/31/11	10/31/11
EPA 8260B	SURROGATE: TOLUENE-D8 (S)	98.9	85-120			%	10/31/11	10/31/11

Quant Method: CALLW.M
Run #: 1030C36
Instrument: Chico
Sequence: C111030
Dilution Factor: 1
Initials: ARS

Printed: 12/06/11 3:06:18 PM
APPL-F1-SC-NoMC-REG MDLs

Quantitation Report (QT Reviewed)

Data File : M:\CHICO\DATA\C111030\1030C36W.D Vial: 1
Acq On : 31 Oct 11 14:16 Operator: STC
Sample : AY49333W04 Inst : Chico
Misc : Water 10mLw/ IS&S:10-30/10-26-11 Multiplr: 1.00

Quant Time: Nov 3 11:55 2011 Quant Results File: CALLW.RES

Quant Method : M:\CHICO\DATA\C111030\CALLW.M (RTE Integrator)
Title : METHOD 8260
Last Update : Wed Nov 02 14:33:25 2011
Response via : Initial Calibration
DataAcq Meth : V8260

Internal Standards	R.T.	QIon	Response	Conc Units		Dev(Min)
1) Fluorobenzene (IS)	12.85	96	558622	25.00000 ppb		0.02
55) Chlorobenzene-D5 (IS)	18.05	117	391168	25.00000 ppb		0.00
71) 1,4-Dichlorobenzene-D (IS)	22.26	152	206336	25.00000 ppb		0.02
System Monitoring Compounds						
33) Dibromofluoromethane(S)	11.43	111	393930	26.47220 ppb		0.00
Spiked Amount 25.097			Recovery = 105.478%			
38) 1,2-DCA-D4(S)	12.23	65	365380	27.58295 ppb		0.00
Spiked Amount 24.225			Recovery = 113.860%			
56) Toluene-D8(S)	15.51	98	1404152	25.51187 ppb		0.02
Spiked Amount 25.808			Recovery = 98.852%			
64) 4-Bromofluorobenzene(S)	20.12	95	515677	26.14866 ppb		0.00
Spiked Amount 25.459			Recovery = 102.708%			
Target Compounds						Qvalue
73) Isopropylbenzene	19.75	105	376343	5.03244 ppb		100
78) n-Propylbenzene	20.46	91	491031	5.50399 ppb		99
83) Tert-Butylbenzene	21.38	119	58606	0.89124 ppb		96
85) Sec-Butylbenzene	21.76	105	425161	5.39136 ppb		96
91) n-Butylbenzene	22.71	91	175843	2.98472 ppb	#	87
96) Naphthalene	25.94	128	360908	14.38582 ppb		95

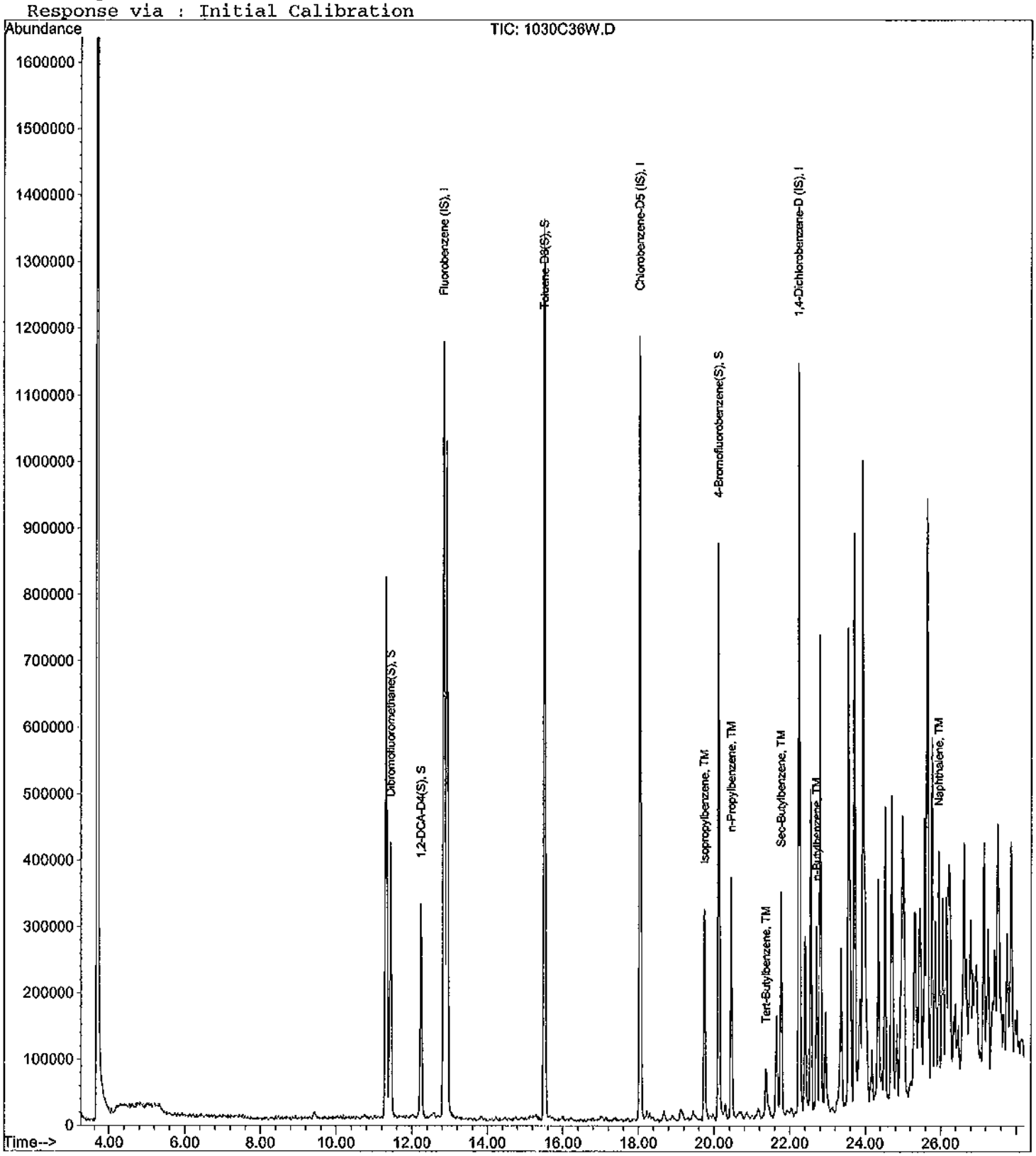
(#) = qualifier out of range (m) = manual integration

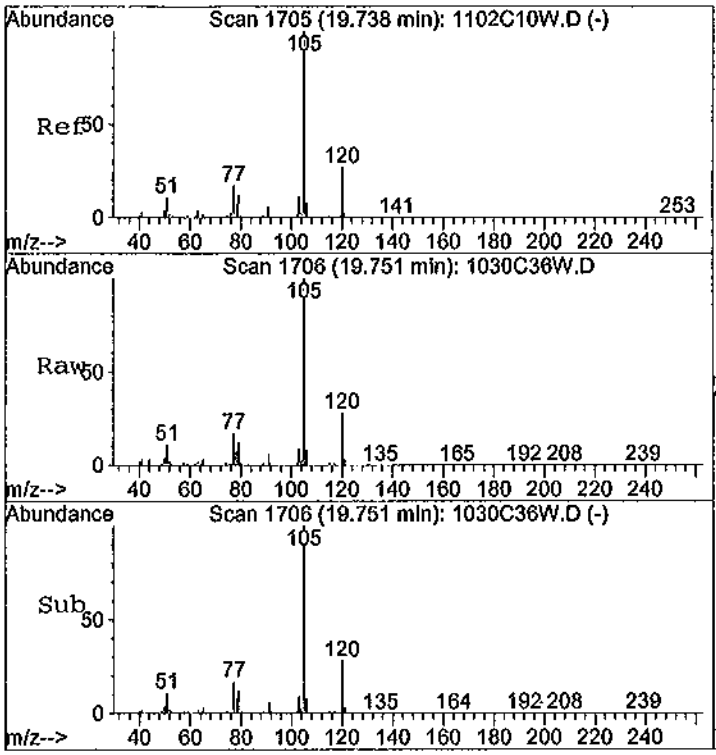
Quantitation Report

Data File : M:\CHICO\DATA\C111030\1030C36W.D Vial: 1
Acq On : 31 Oct 11 14:16 Operator: STC
Sample : AY49333W04 Inst : Chico
Misc : Water 10mLw/ IS&S:10-30/10-26-11 Multiplr: 1.00

Quant Time: Nov 3 11:55 2011 Quant Results File: CALLW.RES

Method : M:\CHICO\DATA\C111030\CALLW.M (RTE Integrator)
Title : METHOD 8260
Last Update : Thu Nov 03 10:27:07 2011
Response via : Initial Calibration

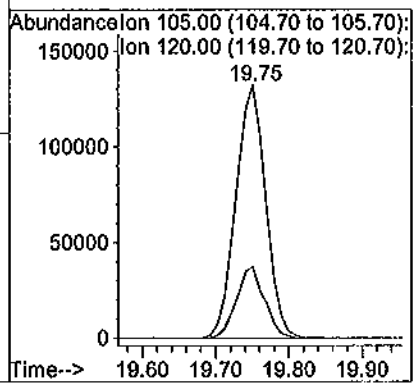


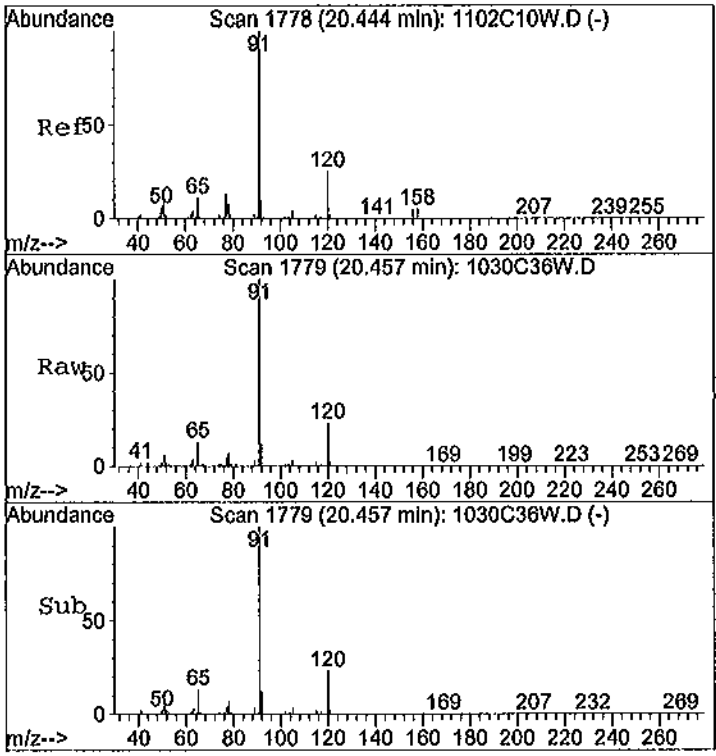


#73
Isopropylbenzene
Concen: 5.03244 ppb
RT: 19.75 min Scan# 1706
Delta R.T. 0.02 min
Lab File: 1030C36W.D
Acq: 31 Oct 11 14:16

Tgt Ion:105 Resp: 376343

Ion	Ratio	Lower	Upper
105	100		
120	28.1	22.6	33.8

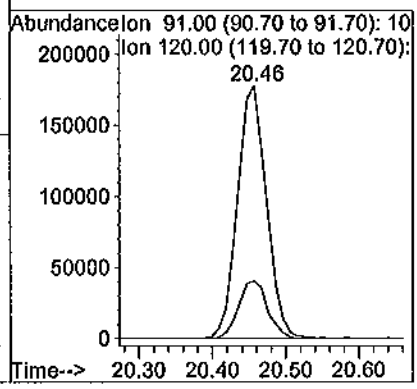


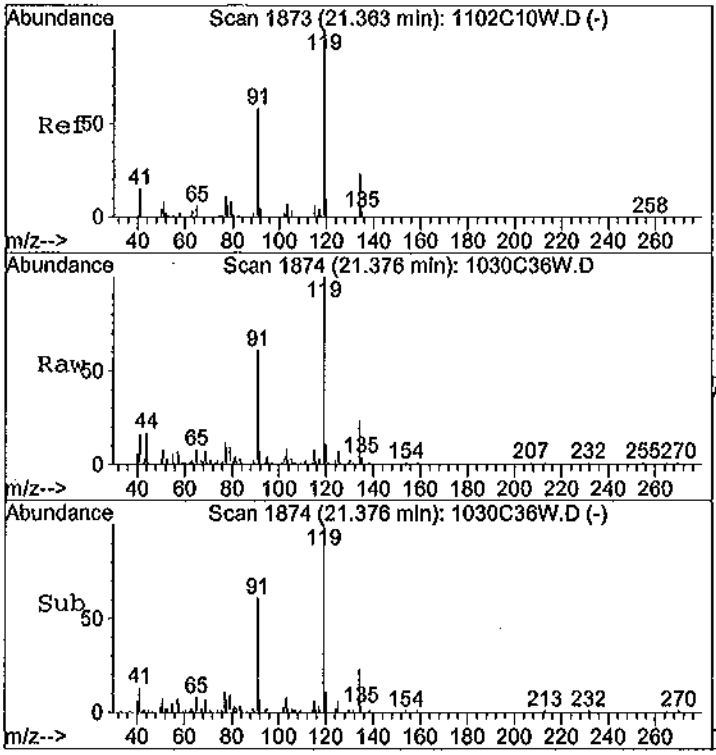


#78
n-Propylbenzene
Concen: 5.50399 ppb
RT: 20.46 min Scan# 1779
Delta R.T. 0.02 min
Lab File: 1030C36W.D
Acq: 31 Oct 11 14:16

Tgt Ion: 91 Resp: 491031

Ion	Ratio	Lower	Upper
91	100		
120	22.8	15.7	29.1

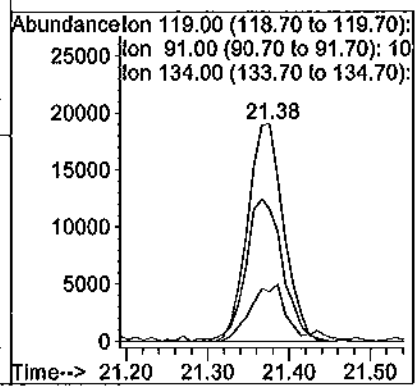


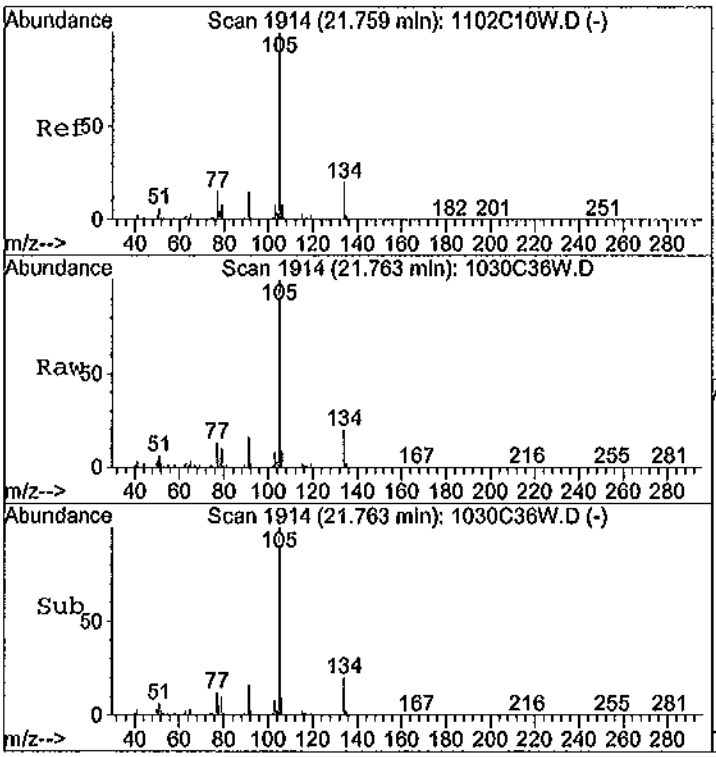


#83
Tert-Butylbenzene
Concen: 0.89124 ppb
RT: 21.38 min Scan# 1874
Delta R.T. 0.02 min
Lab File: 1030C36W.D
Acq: 31 Oct 11 14:16

Tgt Ion:119 Resp: 58606

Ion	Ratio	Lower	Upper
119	100		
91	60.7	43.5	80.7
134	22.8	18.8	34.8

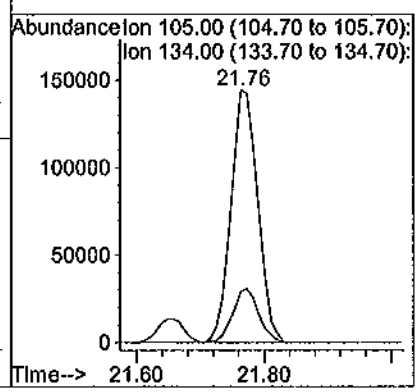


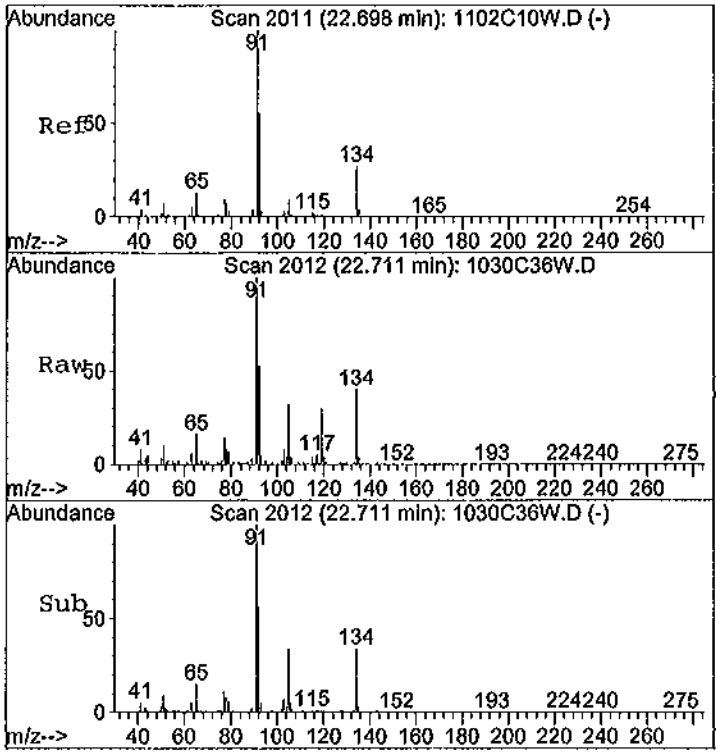


#85
Sec-Butylbenzene
Concen: 5.39136 ppb
RT: 21.76 min Scan# 1914
Delta R.T. 0.01 min
Lab File: 1030C36W.D
Acq: 31 Oct 11 14:16

Tgt Ion:105 Resp: 425161

Ion	Ratio	Lower	Upper
105	100		
134	20.2	15.3	28.5

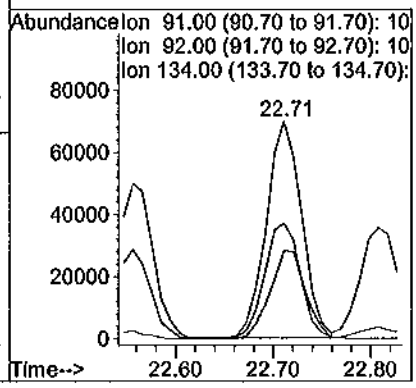


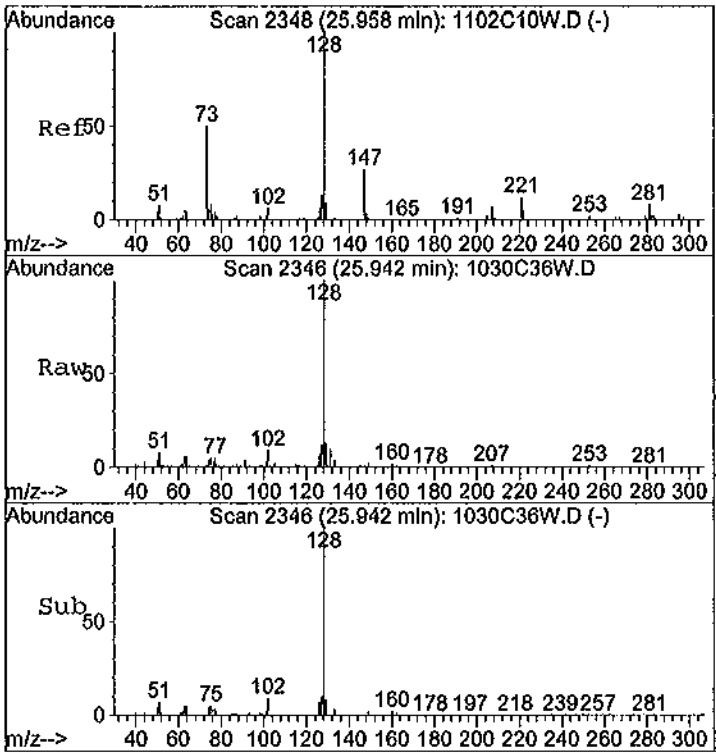


#91
n-Butylbenzene
Concen: 2.98472 ppb
RT: 22.71 min Scan# 2012
Delta R.T. 0.02 min
Lab File: 1030C36W.D
Acq: 31 Oct 11 14:16

Tgt Ion: 91 Resp: 175843

Ion	Ratio	Lower	Upper
91	100		
92	53.2	41.4	77.0
134	40.1	19.8	36.8#

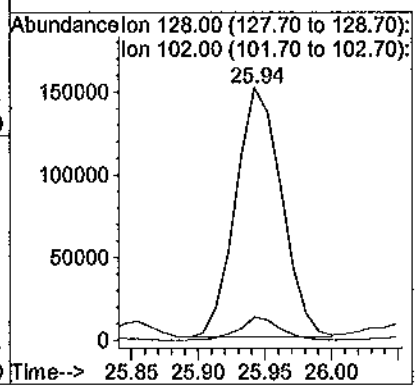




#96
Naphthalene
Concen: 14.38582 ppb
RT: 25.94 min Scan# 2346
Delta R.T. 0.01 min
Lab File: 1030C36W.D
Acq: 31 Oct 11 14:16

Tgt Ion:128 Resp: 360908

Ion	Ratio	Lower	Upper
128	100		
102	9.2	5.1	9.5



Quantitation Report (QT Reviewed)

Data File : M:\CHICO\DATA\C111030\1030C36W.D Vial: 1
Acq On : 31 Oct 11 14:16 Operator: STC
Sample : AY49333W04 Inst : Chico
Misc : Water 10mLw/ IS&S:10-30/10-26-11 Multiplr: 1.00

Quant Time: Nov 10 10:23 2011 Quant Results File: CGAS.RES

Quant Method : M:\CHICO\DATA\C111030\CGAS.M (RTE Integrator)
Title : METHOD 8260B
Last Update : Thu Nov 03 10:47:02 2011
Response via : Initial Calibration
DataAcq Meth : V8260

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Fluorobenzene (IS)	12.85	TIC	1167658	25.00000	ppb	0.01
3) Chlorobenzene-D5 (IS)	18.05	TIC	1180896	25.00000	ppb	0.01
4) 1,4-Dichlorobenzene-D (IS)	22.25	TIC	1128483	25.00000	ppb	0.00

System Monitoring Compounds

Target Compounds Qvalue

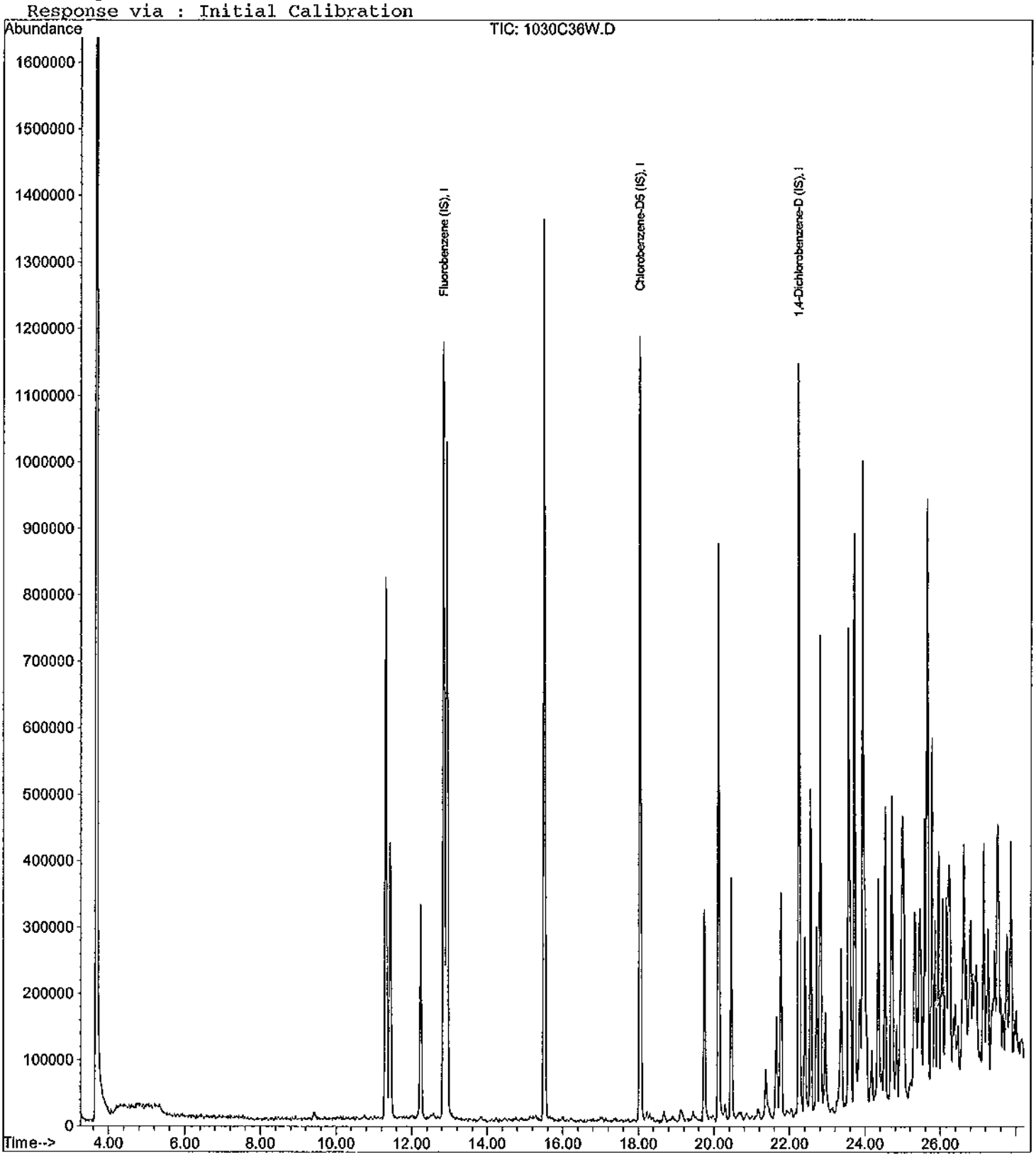
(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : M:\CHICO\DATA\C111030\1030C36W.D Vial: 1
Acq On : 31 Oct 11 14:16 Operator: STC
Sample : AY49333W04 Inst : Chico
Misc : Water 10mLw/ IS&S:10-30/10-26-11 Multiplr: 1.00

Quant Time: Nov 10 10:23 2011 Quant Results File: CGAS.RES

Method : M:\CHICO\DATA\C111030\CGAS.M (RTE Integrator)
Title : METHOD 8260B
Last Update : Thu Nov 03 10:47:02 2011
Response via : Initial Calibration



EPA 8260B VOCs + Gas Water

Environet, Inc.
650 Iwilei Rd, #204
Honolulu, HI 96817

APPL Inc.
908 North Temperance Avenue
Clovis, CA 93611

Attn: Stacy Fineran

Project: RED HILL/1022-024

ARF: 66102

Sample ID: ES047

APPL ID: AY49334

Sample Collection Date: 10/24/11

QCG: #86RHB-111030AC-161029

Method	Analyte	Result	LOQ	LOD	DL	Units	Extraction Date	Analysis Date
EPA 8260B	1,1,1,2-TETRACHLOROETHANE	0.26 U	1.0	0.26	0.13	ug/L	10/31/11	10/31/11
EPA 8260B	1,1,1-TRICHLOROETHANE	0.28 U	1.0	0.28	0.14	ug/L	10/31/11	10/31/11
EPA 8260B	1,1,2,2-TETRACHLOROETHANE	0.20 U	1.0	0.20	0.10	ug/L	10/31/11	10/31/11
EPA 8260B	1,1,2-TRICHLOROETHANE	0.40 U	1.0	0.40	0.20	ug/L	10/31/11	10/31/11
EPA 8260B	1,1-DICHLOROETHANE	0.38 U	1.0	0.38	0.19	ug/L	10/31/11	10/31/11
EPA 8260B	1,1-DICHLOROETHENE	0.60 U	1.0	0.60	0.30	ug/L	10/31/11	10/31/11
EPA 8260B	1,2,3-TRICHLOROPROPANE	0.78 U	2.0	0.78	0.39	ug/L	10/31/11	10/31/11
EPA 8260B	1,2,4-TRICHLOROBENZENE	0.42 U	1.0	0.42	0.21	ug/L	10/31/11	10/31/11
EPA 8260B	1,2-DIBROMO-3-CHLOROPROPANE	1.52 U	2.0	1.52	0.76	ug/L	10/31/11	10/31/11
EPA 8260B	1,2-DIBROMOETHANE	0.40 U	1.0	0.40	0.20	ug/L	10/31/11	10/31/11
EPA 8260B	1,2-DICHLOROBENZENE	0.34 U	1.0	0.34	0.17	ug/L	10/31/11	10/31/11
EPA 8260B	1,2-DICHLOROETHANE	0.28 U	1.0	0.28	0.14	ug/L	10/31/11	10/31/11
EPA 8260B	1,2-DICHLOROPROPANE	0.34 U	1.0	0.34	0.17	ug/L	10/31/11	10/31/11
EPA 8260B	1,3-DICHLOROBENZENE	0.22 U	1.0	0.22	0.11	ug/L	10/31/11	10/31/11
EPA 8260B	1,3-DICHLOROPROPENE, TOTAL	0.36 U	1.0	0.36	0.18	ug/L	10/31/11	10/31/11
EPA 8260B	1,4-DICHLOROBENZENE	0.38 U	1.0	0.38	0.19	ug/L	10/31/11	10/31/11
EPA 8260B	2-BUTANONE	1.20 U	10.0	1.20	0.60	ug/L	10/31/11	10/31/11
EPA 8260B	4-METHYL-2-PENTANONE	3.80 U	10.0	3.80	1.90	ug/L	10/31/11	10/31/11
EPA 8260B	ACETONE	1.90 U	10.0	1.90	0.95	ug/L	10/31/11	10/31/11
EPA 8260B	BENZENE	0.32 U	1.0	0.32	0.16	ug/L	10/31/11	10/31/11
EPA 8260B	BROMODICHLOROMETHANE	0.28 U	1.0	0.28	0.14	ug/L	10/31/11	10/31/11
EPA 8260B	BROMOFORM	0.28 U	1.0	0.28	0.14	ug/L	10/31/11	10/31/11
EPA 8260B	BROMOMETHANE	0.48 U	2.0	0.48	0.24	ug/L	10/31/11	10/31/11
EPA 8260B	CARBON TETRACHLORIDE	0.20 U	1.0	0.20	0.10	ug/L	10/31/11	10/31/11
EPA 8260B	CHLOROBENZENE	0.42 U	1.0	0.42	0.21	ug/L	10/31/11	10/31/11
EPA 8260B	CHLORODIBROMOMETHANE	0.38 U	1.0	0.38	0.19	ug/L	10/31/11	10/31/11
EPA 8260B	CHLOROETHANE	0.42 U	1.0	0.42	0.21	ug/L	10/31/11	10/31/11
EPA 8260B	CHLOROFORM	0.14 U	1.0	0.14	0.07	ug/L	10/31/11	10/31/11
EPA 8260B	CHLOROMETHANE	0.84 U	1.0	0.84	0.42	ug/L	10/31/11	10/31/11
EPA 8260B	CIS-1,2-DICHLOROETHENE	0.32 U	1.0	0.32	0.16	ug/L	10/31/11	10/31/11
EPA 8260B	ETHYLBENZENE	0.46 U	1.0	0.46	0.23	ug/L	10/31/11	10/31/11
EPA 8260B	GASOLINE	12.12 U	20.0	12.12	6.06	ug/L	10/31/11	10/31/11
EPA 8260B	HEXACHLOROBUTADIENE	0.38 U	1.0	0.38	0.19	ug/L	10/31/11	10/31/11
EPA 8260B	METHYL TERT-BUTYL ETHER	0.52 U	1.0	0.52	0.26	ug/L	10/31/11	10/31/11
EPA 8260B	METHYLENE CHLORIDE	0.70 U	5.0	0.70	0.35	ug/L	10/31/11	10/31/11

Quant Method: CALLW.M
Run #: 1030C37
Instrument: Chico
Sequence: C111030
Dilution Factor: 1
Initials: ARS

Printed: 12/06/11 3:06:18 PM

APPL-F1-SC-NoMC-REG MDLs

EPA 8260B VOCs + Gas Water

Environet, Inc.
650 Iwilei Rd, #204
Honolulu, HI 96817

APPL Inc.
908 North Temperance Avenue
Clovis, CA 93611

Attn: Stacy Fineran

Project: RED HILL/1022-024

ARF: 66102

Sample ID: ES047

APPL ID: AY49334

Sample Collection Date: 10/24/11

QCG: #86RHB-111030AC-161029

Method	Analyte	Result	LOQ	LOD	DL	Units	Extraction Date	Analysis Date
EPA 8260B	STYRENE	0.50 U	1.0	0.50	0.25	ug/L	10/31/11	10/31/11
EPA 8260B	TETRACHLOROETHENE	0.48 U	1.0	0.48	0.24	ug/L	10/31/11	10/31/11
EPA 8260B	TOLUENE	0.34 U	1.0	0.34	0.17	ug/L	10/31/11	10/31/11
EPA 8260B	TRANS-1,2-DICHLOROETHENE	0.38 U	1.0	0.38	0.19	ug/L	10/31/11	10/31/11
EPA 8260B	TRICHLOROETHENE	0.32 U	1.0	0.32	0.16	ug/L	10/31/11	10/31/11
EPA 8260B	VINYL CHLORIDE	0.46 U	1.0	0.46	0.23	ug/L	10/31/11	10/31/11
EPA 8260B	XYLENES (TOTAL)	0.38 U	1.0	0.38	0.19	ug/L	10/31/11	10/31/11
EPA 8260B	SURROGATE: 1,2-DICHLOROETHANE-	104	70-120			%	10/31/11	10/31/11
EPA 8260B	SURROGATE: 4-BROMOFLUOROBEN	98.9	75-120			%	10/31/11	10/31/11
EPA 8260B	SURROGATE: DIBROMOFLUOROMET	102	85-115			%	10/31/11	10/31/11
EPA 8260B	SURROGATE: TOLUENE-D8 (S)	97.6	85-120			%	10/31/11	10/31/11

Quant Method: CALLW.M
Run #: 1030C37
Instrument: Chico
Sequence: C111030
Dilution Factor: 1
Initials: ARS

Printed: 12/06/11 3:06:18 PM
APPL-F1-SC-NoMC-REG MDLs

Quantitation Report (QT Reviewed)

Data File : M:\CHICO\DATA\C111030\1030C37W.D Vial: 1
Acq On : 31 Oct 11 14:53 Operator: STC
Sample : AY49334W13 Inst : Chico
Misc : Water 10mLw/ IS&S:10-30/10-26-11 Multiplr: 1.00

Quant Time: Nov 3 12:00 2011 Quant Results File: CALLW.RES

Quant Method : M:\CHICO\DATA\C111030\CALLW.M (RTE Integrator)
Title : METHOD 8260
Last Update : Wed Nov 02 14:33:25 2011
Response via : Initial Calibration
DataAcq Meth : V8260

Internal Standards	R.T.	QIon	Response	Conc Units	Dev(Min)
1) Fluorobenzene (IS)	12.85	96	617536	25.00000 ppb	0.02
55) Chlorobenzene-D5 (IS)	18.05	117	437056	25.00000 ppb	0.00
71) 1,4-Dichlorobenzene-D (IS)	22.26	152	232960	25.00000 ppb	0.02
System Monitoring Compounds					
33) Dibromofluoromethane(S)	11.43	111	420841	25.58260 ppb	0.00
Spiked Amount 25.097			Recovery	= 101.935%	
38) 1,2-DCA-D4(S)	12.23	65	368967	25.19644 ppb	0.00
Spiked Amount 24.225			Recovery	= 104.007%	
56) Toluene-D8(S)	15.51	98	1549428	25.19567 ppb	0.02
Spiked Amount 25.808			Recovery	= 97.628%	
64) 4-Bromofluorobenzene(S)	20.13	95	554811	25.17926 ppb	0.02
Spiked Amount 25.459			Recovery	= 98.898%	
Target Compounds					Qvalue
73) Isopropylbenzene	19.75	105	388062	4.59610 ppb	99
78) n-Propylbenzene	20.46	91	482750	4.79275 ppb	100
83) Tert-Butylbenzene	21.37	119	65995	0.88891 ppb	94
85) Sec-Butylbenzene	21.77	105	453460	5.09305 ppb	97
96) Naphthalene	25.94	128	258129	9.11315 ppb	95

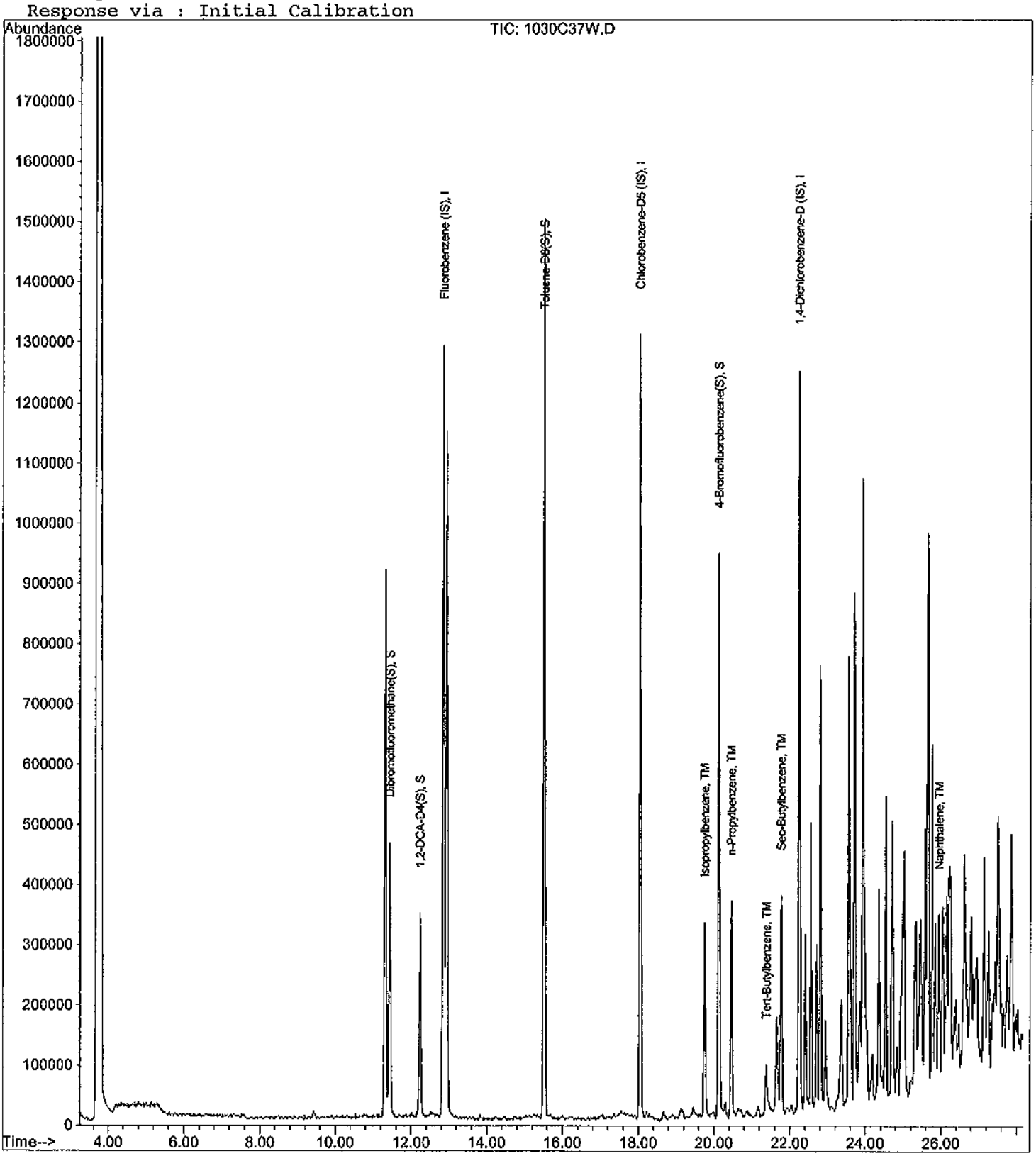
(#) = qualifier out of range (m) = manual integration

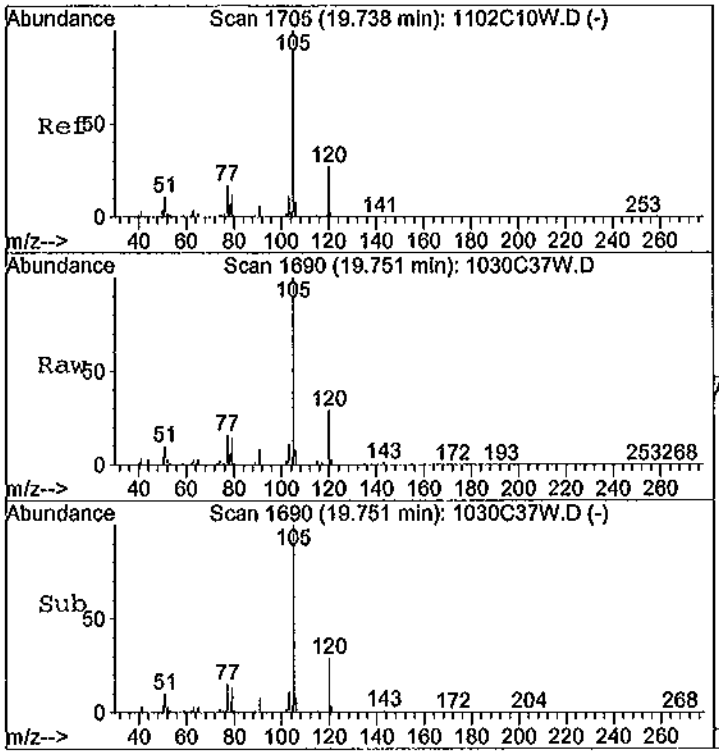
Quantitation Report

Data File : M:\CHICO\DATA\C111030\1030C37W.D Vial: 1
Acq On : 31 Oct 11 14:53 Operator: STC
Sample : AY49334W13 Inst : Chico
Misc : Water 10mLw/ IS&S:10-30/10-26-11 Multiplr: 1.00

Quant Time: Nov 3 12:00 2011 Quant Results File: CALLW.RES

Method : M:\CHICO\DATA\C111030\CALLW.M (RTE Integrator)
Title : METHOD 8260
Last Update : Thu Nov 03 10:27:07 2011
Response via : Initial Calibration

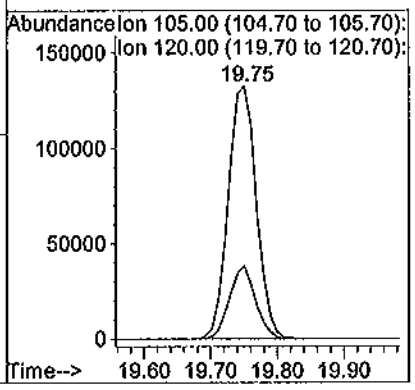


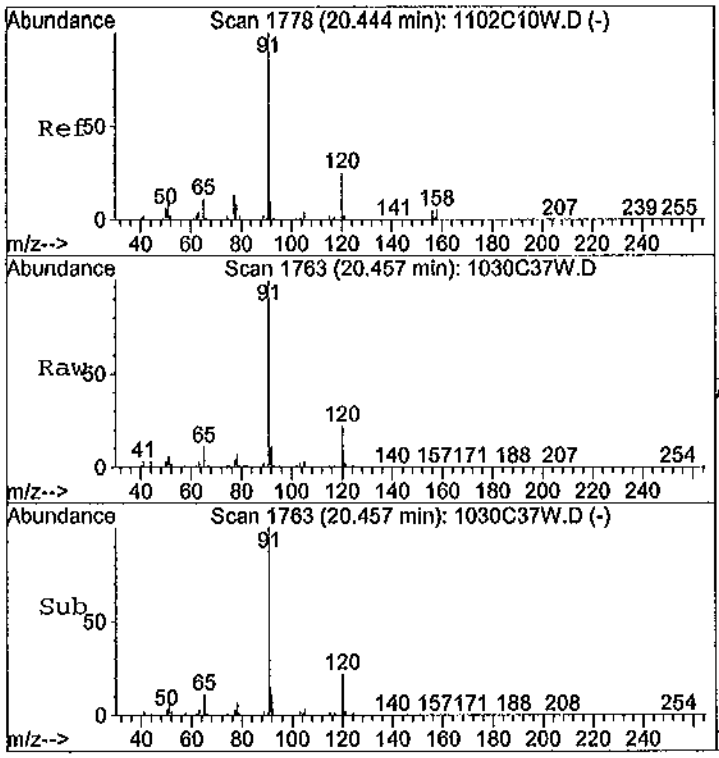


#73
Isopropylbenzene
Concen: 4.59610 ppb
RT: 19.75 min Scan# 1690
Delta R.T. 0.02 min
Lab File: 1030C37W.D
Acq: 31 Oct 11 14:53

Tgt Ion:105 Resp: 388062

Ion	Ratio	Lower	Upper
105	100		
120	28.8	22.6	33.8

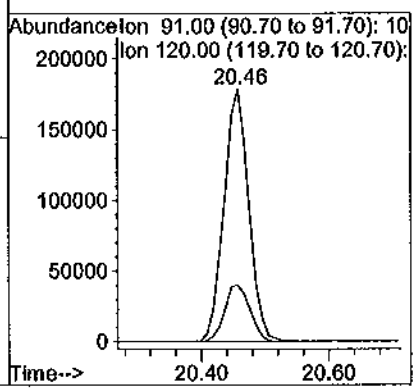


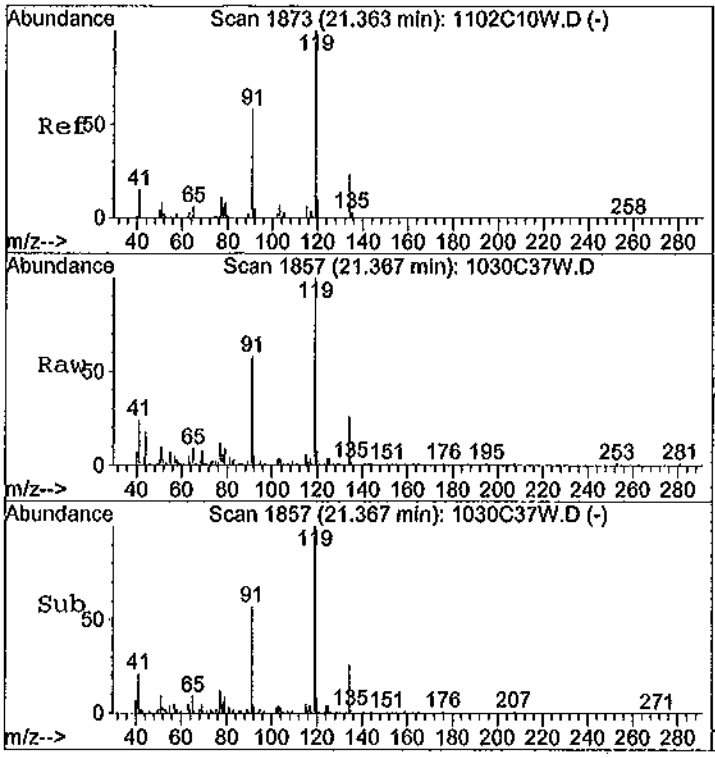


#78
n-Propylbenzene
Concen: 4.79275 ppb
RT: 20.46 min Scan# 1763
Delta R.T. 0.02 min
Lab File: 1030C37W.D
Acq: 31 Oct 11 14:53

Tgt Ion: 91 Resp: 482750

Ion	Ratio	Lower	Upper
91	100		
120	22.3	15.7	29.1

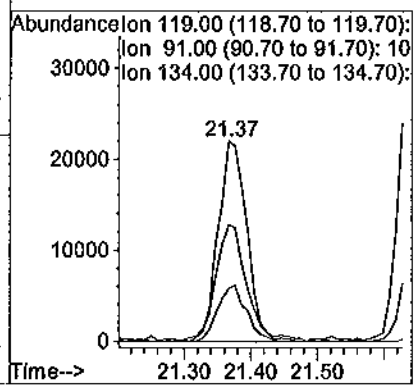


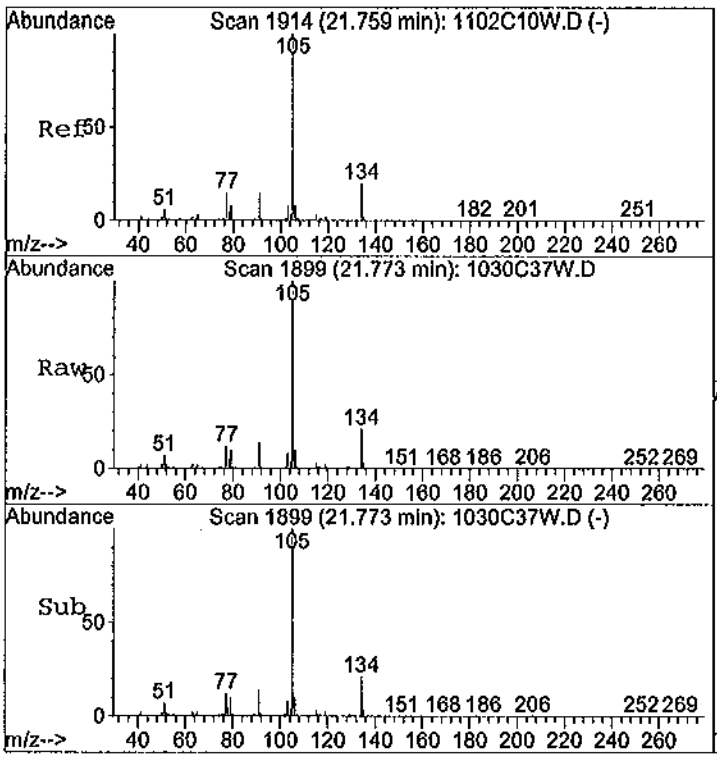


#83
Tert-Butylbenzene
Concen: 0.88891 ppb
RT: 21.37 min Scan# 1857
Delta R.T. 0.01 min
Lab File: 1030C37W.D
Acq: 31 Oct 11 14:53

Tgt Ion:119 Resp: 65995

Ion	Ratio	Lower	Upper
119	100		
91	56.1	43.5	80.7
134	26.3	18.8	34.8

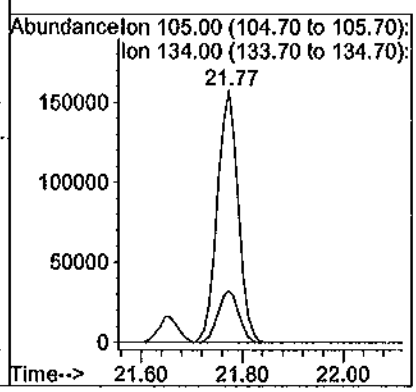


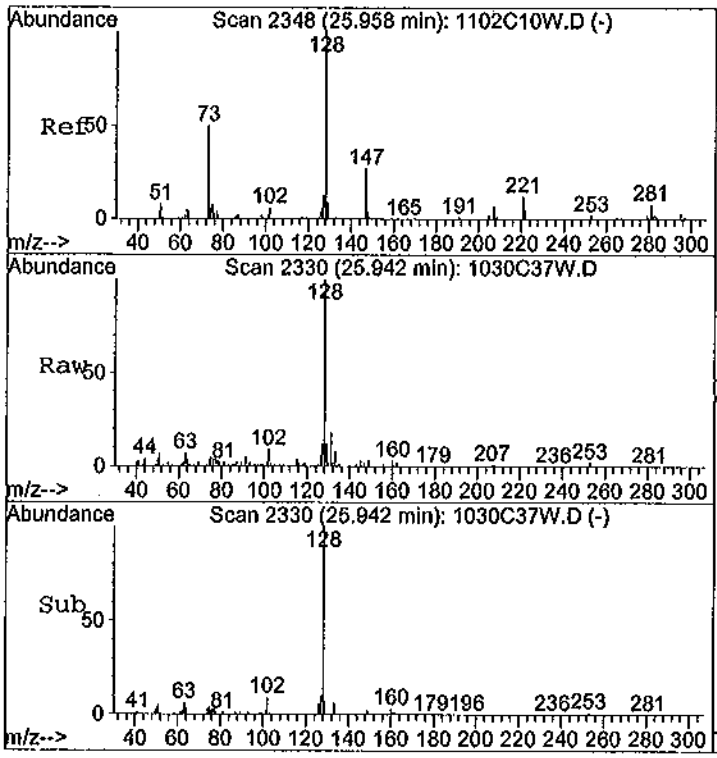


#85
Sec-Butylbenzene
Concen: 5.09305 ppb
RT: 21.77 min Scan# 1899
Delta R.T. 0.02 min
Lab File: 1030C37W.D
Acq: 31 Oct 11 14:53

Tgt Ion:105 Resp: 453460

Ion	Ratio	Lower	Upper
105	100		
134	20.6	15.3	28.5

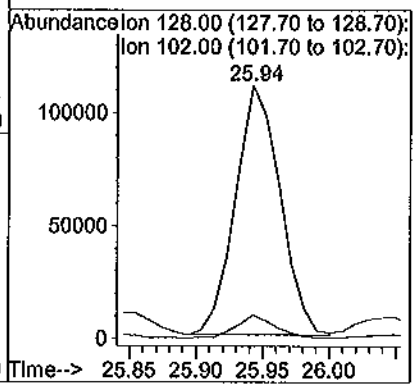




#96
Naphthalene
Concen: 9.11315 ppb
RT: 25.94 min Scan# 2330
Delta R.T. 0.01 min
Lab File: 1030C37W.D
Acq: 31 Oct 11 14:53

Tgt Ion:128 Resp: 258129

Ion	Ratio	Lower	Upper
128	100		
102	9.2	5.1	9.5



Quantitation Report (QT Reviewed)

Data File : M:\CHICO\DATA\C111030\1030C37W.D Vial: 1
Acq On : 31 Oct 11 14:53 Operator: STC
Sample : AY49334W13 Inst : Chico
Misc : Water 10mLw/ IS&S:10-30/10-26-11 Multiplr: 1.00

Quant Time: Nov 10 10:23 2011 Quant Results File: CGAS.RES

Quant Method : M:\CHICO\DATA\C111030\CGAS.M (RTE Integrator)
Title : METHOD 8260B
Last Update : Thu Nov 03 10:47:02 2011
Response via : Initial Calibration
DataAcq Meth : V8260

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Fluorobenzene (IS)	12.85	TIC	1282349	25.00000	ppb	0.01
3) Chlorobenzene-D5 (IS)	18.05	TIC	1302320	25.00000	ppb	0.01
4) 1,4-Dichlorobenzene-D (IS)	22.26	TIC	1229710	25.00000	ppb	0.01

System Monitoring Compounds

Target Compounds Qvalue

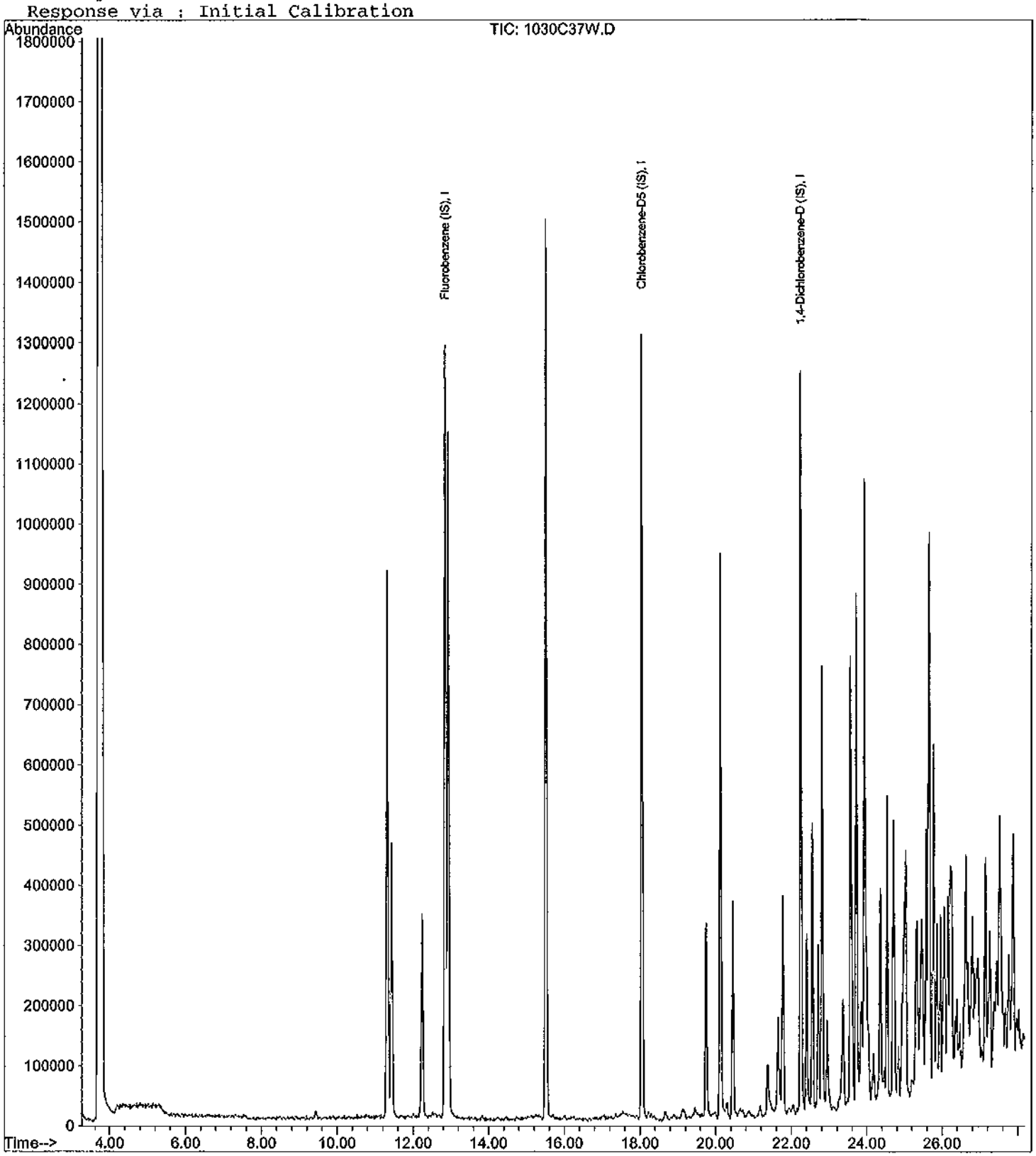
(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : M:\CHICO\DATA\C111030\1030C37W.D Vial: 1
Acq On : 31 Oct 11 14:53 Operator: STC
Sample : AY49334W13 Inst : Chico
Misc : Water 10mLw/ IS&S:10-30/10-26-11 Multiplr: 1.00

Quant Time: Nov 10 10:23 2011 Quant Results File: CGAS.RES

Method : M:\CHICO\DATA\C111030\CGAS.M (RTE Integrator)
Title : METHOD 8260B
Last Update : Thu Nov 03 10:47:02 2011
Response via : Initial Calibration



EPA 8260B VOCs + Gas Water

Environet, Inc.
650 Iwilei Rd, #204
Honolulu, HI 96817

Attn: Stacy Fineran
Project: RED HILL/1022-024
Sample ID: ES048
Sample Collection Date: 10/24/11

APPL Inc.
908 North Temperance Avenue
Clovis, CA 93611

ARF: 66102
APPL ID: AY49335
QCG: #86RHB-111030AC-161029

Method	Analyte	Result	LOQ	LOD	DL	Units	Extraction Date	Analysis Date
EPA 8260B	1,1,1,2-TETRACHLOROETHANE	0.26 U	1.0	0.26	0.13	ug/L	10/31/11	10/31/11
EPA 8260B	1,1,1-TRICHLOROETHANE	0.28 U	1.0	0.28	0.14	ug/L	10/31/11	10/31/11
EPA 8260B	1,1,2,2-TETRACHLOROETHANE	0.20 U	1.0	0.20	0.10	ug/L	10/31/11	10/31/11
EPA 8260B	1,1,2-TRICHLOROETHANE	0.40 U	1.0	0.40	0.20	ug/L	10/31/11	10/31/11
EPA 8260B	1,1-DICHLOROETHANE	0.38 U	1.0	0.38	0.19	ug/L	10/31/11	10/31/11
EPA 8260B	1,1-DICHLOROETHENE	0.60 U	1.0	0.60	0.30	ug/L	10/31/11	10/31/11
EPA 8260B	1,2,3-TRICHLOROPROPANE	0.78 U	2.0	0.78	0.39	ug/L	10/31/11	10/31/11
EPA 8260B	1,2,4-TRICHLOROBENZENE	0.42 U	1.0	0.42	0.21	ug/L	10/31/11	10/31/11
EPA 8260B	1,2-DIBROMO-3-CHLOROPROPANE	1.52 U	2.0	1.52	0.76	ug/L	10/31/11	10/31/11
EPA 8260B	1,2-DIBROMOETHANE	0.40 U	1.0	0.40	0.20	ug/L	10/31/11	10/31/11
EPA 8260B	1,2-DICHLOROBENZENE	0.34 U	1.0	0.34	0.17	ug/L	10/31/11	10/31/11
EPA 8260B	1,2-DICHLOROETHANE	0.28 U	1.0	0.28	0.14	ug/L	10/31/11	10/31/11
EPA 8260B	1,2-DICHLOROPROPANE	0.34 U	1.0	0.34	0.17	ug/L	10/31/11	10/31/11
EPA 8260B	1,3-DICHLOROBENZENE	0.22 U	1.0	0.22	0.11	ug/L	10/31/11	10/31/11
EPA 8260B	1,3-DICHLOROPROPENE, TOTAL	0.36 U	1.0	0.36	0.18	ug/L	10/31/11	10/31/11
EPA 8260B	1,4-DICHLOROBENZENE	0.38 U	1.0	0.38	0.19	ug/L	10/31/11	10/31/11
EPA 8260B	2-BUTANONE	1.20 U	10.0	1.20	0.60	ug/L	10/31/11	10/31/11
EPA 8260B	4-METHYL-2-PENTANONE	3.80 U	10.0	3.80	1.90	ug/L	10/31/11	10/31/11
EPA 8260B	ACETONE	1.90 U	10.0	1.90	0.95	ug/L	10/31/11	10/31/11
EPA 8260B	BENZENE	0.32 U	1.0	0.32	0.16	ug/L	10/31/11	10/31/11
EPA 8260B	BROMODICHLOROMETHANE	0.28 U	1.0	0.28	0.14	ug/L	10/31/11	10/31/11
EPA 8260B	BROMOFORM	0.28 U	1.0	0.28	0.14	ug/L	10/31/11	10/31/11
EPA 8260B	BROMOMETHANE	0.48 U	2.0	0.48	0.24	ug/L	10/31/11	10/31/11
EPA 8260B	CARBON TETRACHLORIDE	0.20 U	1.0	0.20	0.10	ug/L	10/31/11	10/31/11
EPA 8260B	CHLOROBENZENE	0.42 U	1.0	0.42	0.21	ug/L	10/31/11	10/31/11
EPA 8260B	CHLORODIBROMOMETHANE	0.38 U	1.0	0.38	0.19	ug/L	10/31/11	10/31/11
EPA 8260B	CHLOROETHANE	0.42 U	1.0	0.42	0.21	ug/L	10/31/11	10/31/11
EPA 8260B	CHLOROFORM	0.14 U	1.0	0.14	0.07	ug/L	10/31/11	10/31/11
EPA 8260B	CHLOROMETHANE	0.84 U	1.0	0.84	0.42	ug/L	10/31/11	10/31/11
EPA 8260B	CIS-1,2-DICHLOROETHENE	0.32 U	1.0	0.32	0.16	ug/L	10/31/11	10/31/11
EPA 8260B	ETHYLBENZENE	0.46 U	1.0	0.46	0.23	ug/L	10/31/11	10/31/11
EPA 8260B	GASOLINE	12.12 U	20.0	12.12	6.06	ug/L	10/31/11	10/31/11
EPA 8260B	HEXACHLOROBUTADIENE	0.38 U	1.0	0.38	0.19	ug/L	10/31/11	10/31/11
EPA 8260B	METHYL TERT-BUTYL ETHER	0.52 U	1.0	0.52	0.26	ug/L	10/31/11	10/31/11
EPA 8260B	METHYLENE CHLORIDE	0.70 U	5.0	0.70	0.35	ug/L	10/31/11	10/31/11

Quant Method: CALLW.M
Run #: 1030C35
Instrument: Chico
Sequence: C111030
Dilution Factor: 1
Initials: ARS

Printed: 12/06/11 3:06:18 PM
APPL-F1-SC-NoMC-REG MDLs

EPA 8260B VOCs + Gas Water

Environet, Inc.
650 Iwilei Rd, #204
Honolulu, HI 96817

APPL Inc.
908 North Temperance Avenue
Clovis, CA 93611

Attn: Stacy Fineran

Project: RED HILL/1022-024

ARF: 66102

Sample ID: ES048

APPL ID: AY49335

Sample Collection Date: 10/24/11

QCG: #86RHB-111030AC-161029

Method	Analyte	Result	LOQ	LOD	DL	Units	Extraction Date	Analysis Date
EPA 8260B	STYRENE	0.50 U	1.0	0.50	0.25	ug/L	10/31/11	10/31/11
EPA 8260B	TETRACHLOROETHENE	0.48 U	1.0	0.48	0.24	ug/L	10/31/11	10/31/11
EPA 8260B	TOLUENE	0.34 U	1.0	0.34	0.17	ug/L	10/31/11	10/31/11
EPA 8260B	TRANS-1,2-DICHLOROETHENE	0.38 U	1.0	0.38	0.19	ug/L	10/31/11	10/31/11
EPA 8260B	TRICHLOROETHENE	0.32 U	1.0	0.32	0.16	ug/L	10/31/11	10/31/11
EPA 8260B	VINYL CHLORIDE	0.46 U	1.0	0.46	0.23	ug/L	10/31/11	10/31/11
EPA 8260B	XYLENES (TOTAL)	0.38 U	1.0	0.38	0.19	ug/L	10/31/11	10/31/11
EPA 8260B	SURROGATE: 1,2-DICHLOROETHANE-	108	70-120			%	10/31/11	10/31/11
EPA 8260B	SURROGATE: 4-BROMOFLUOROBEN	102	75-120			%	10/31/11	10/31/11
EPA 8260B	SURROGATE: DIBROMOFLUOROMET	105	85-115			%	10/31/11	10/31/11
EPA 8260B	SURROGATE: TOLUENE-D8 (S)	99.0	85-120			%	10/31/11	10/31/11

Quant Method: CALLW.M
Run #: 1030C35
Instrument: Chico
Sequence: C111030
Dilution Factor: 1
Initials: ARS

Printed: 12/06/11 3:06:18 PM
APPL-F1-SC-NoMC-REG MDLs

Quantitation Report (QT Reviewed)

Data File : M:\CHICO\DATA\C111030\1030C35W.D Vial: 1
Acq On : 31 Oct 11 13:39 Operator: STC
Sample : AY49335W01 Inst : Chico
Misc : Water 10mLw/ IS&S:10-30/10-26-11 Multiplr: 1.00

Quant Time: Nov 3 11:52 2011 Quant Results File: CALLW.RES

Quant Method : M:\CHICO\DATA\C111030\CALLW.M (RTE Integrator)
Title : METHOD 8260
Last Update : Wed Nov 02 14:33:25 2011
Response via : Initial Calibration
DataAcq Meth : V8260

Internal Standards	R.T.	QIon	Response	Conc Units	Dev(Min)
1) Fluorobenzene (IS)	12.85	96	576569	25.00000 ppb	0.01
55) Chlorobenzene-D5 (IS)	18.05	117	396544	25.00000 ppb	0.01
71) 1,4-Dichlorobenzene-D (IS)	22.25	152	214592	25.00000 ppb	0.01
System Monitoring Compounds					
33) Dibromofluoromethane(S)	11.43	111	402959	26.23605 ppb	0.00
Spiked Amount 25.097			Recovery = 104.537%		
38) 1,2-DCA-D4(S)	12.23	65	357575	26.15350 ppb	0.00
Spiked Amount 24.225			Recovery = 107.958%		
56) Toluene-D8(S)	15.51	98	1425024	25.54008 ppb	0.01
Spiked Amount 25.808			Recovery = 98.960%		
64) 4-Bromofluorobenzene(S)	20.12	95	520964	26.05861 ppb	0.01
Spiked Amount 25.459			Recovery = 102.355%		

Target Compounds Qvalue

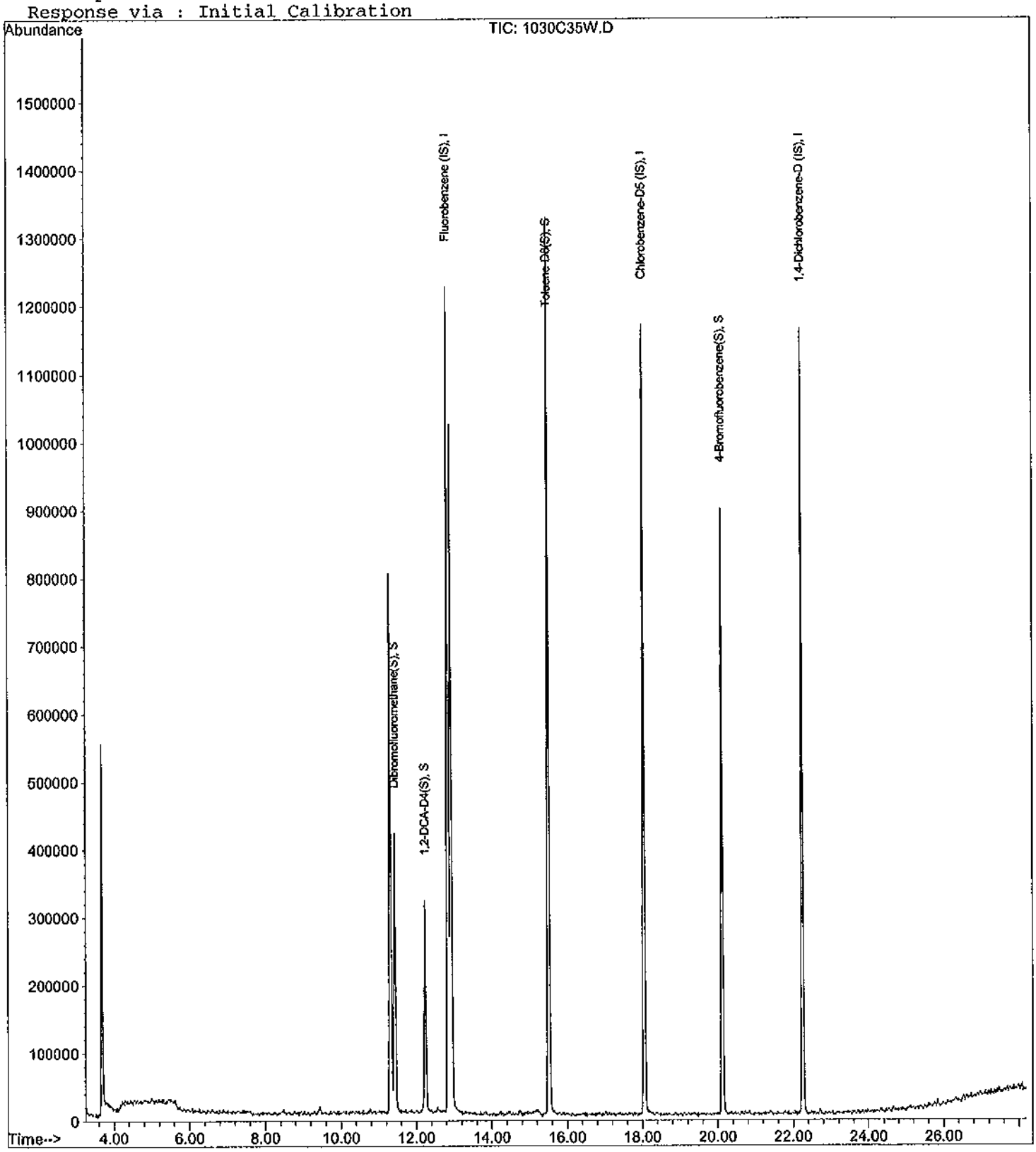
(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : M:\CHICO\DATA\C111030\1030C35W.D Vial: 1
Acq On : 31 Oct 11 13:39 Operator: STC
Sample : AY49335W01 Inst : Chico
Misc : Water 10mLw/ IS&S:10-30/10-26-11 Multiplr: 1.00

Quant Time: Nov 3 11:52 2011 Quant Results File: CALLW.RES

Method : M:\CHICO\DATA\C111030\CALLW.M (RTE Integrator)
Title : METHOD 8260
Last Update : Thu Nov 03 10:27:07 2011
Response via : Initial Calibration



Quantitation Report (QT Reviewed)

Data File : M:\CHICO\DATA\C111030\1030C35W.D Vial: 1
Acq On : 31 Oct 11 13:39 Operator: STC
Sample : AY49335W01 Inst : Chico
Misc : Water 10mLw/ IS&S:10-30/10-26-11 Multiplr: 1.00

Quant Time: Nov 10 10:23 2011 Quant Results File: CGAS.RES

Quant Method : M:\CHICO\DATA\C111030\CGAS.M (RTE Integrator)
Title : METHOD 8260B
Last Update : Thu Nov 03 10:47:02 2011
Response via : Initial Calibration
DataAcq Meth : V8260

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Fluorobenzene (IS)	12.85	TIC	1215846	25.00000	ppb	0.00
3) Chlorobenzene-D5 (IS)	18.05	TIC	1162609	25.00000	ppb	0.02
4) 1,4-Dichlorobenzene-D (IS)	22.25	TIC	1158136	25.00000	ppb	0.00

System Monitoring Compounds

Target Compounds Qvalue

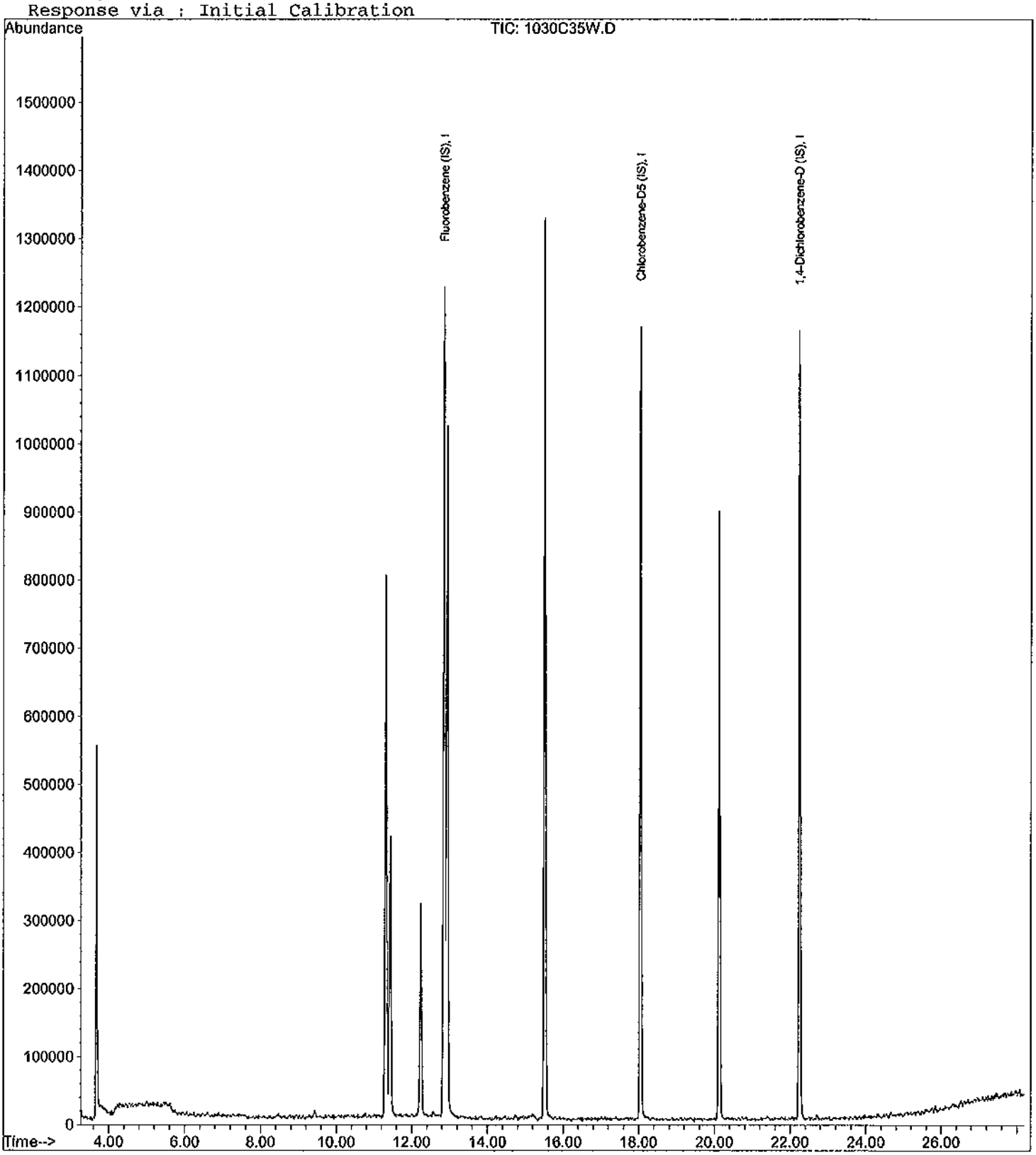
(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : M:\CHICO\DATA\C111030\1030C35W.D Vial: 1
Acq On : 31 Oct 11 13:39 Operator: STC
Sample : AY49335W01 Inst : Chico
Misc : Water 10mLw/ IS&S:10-30/10-26-11 Multiplr: 1.00

Quant Time: Nov 10 10:23 2011 Quant Results File: CGAS.RES

Method : M:\CHICO\DATA\C111030\CGAS.M (RTE Integrator)
Title : METHOD 8260B
Last Update : Thu Nov 03 10:47:02 2011
Response via : Initial Calibration



EPA 8260B VOCs + Gas Water

Environet, Inc.
650 Iwilei Rd, #204
Honolulu, HI 96817

APPL Inc.
908 North Temperance Avenue
Clovis, CA 93611

Attn: Stacy Fineran
Project: RED HILL/1022-024
Sample ID: ES049
Sample Collection Date: 10/24/11

ARF: 66102
APPL ID: AY49336
QCG: #86RHB-111030AC-161029

Method	Analyte	Result	LOQ	LOD	DL	Units	Extraction Date	Analysis Date
EPA 8260B	1,1,1,2-TETRACHLOROETHANE	0.26 U	1.0	0.26	0.13	ug/L	10/31/11	10/31/11
EPA 8260B	1,1,1-TRICHLOROETHANE	0.28 U	1.0	0.28	0.14	ug/L	10/31/11	10/31/11
EPA 8260B	1,1,2,2-TETRACHLOROETHANE	0.20 U	1.0	0.20	0.10	ug/L	10/31/11	10/31/11
EPA 8260B	1,1,2-TRICHLOROETHANE	0.40 U	1.0	0.40	0.20	ug/L	10/31/11	10/31/11
EPA 8260B	1,1-DICHLOROETHANE	0.38 U	1.0	0.38	0.19	ug/L	10/31/11	10/31/11
EPA 8260B	1,1-DICHLOROETHENE	0.60 U	1.0	0.60	0.30	ug/L	10/31/11	10/31/11
EPA 8260B	1,2,3-TRICHLOROPROPANE	0.78 U	2.0	0.78	0.39	ug/L	10/31/11	10/31/11
EPA 8260B	1,2,4-TRICHLOROBENZENE	0.42 U	1.0	0.42	0.21	ug/L	10/31/11	10/31/11
EPA 8260B	1,2-DIBROMO-3-CHLOROPROPANE	1.52 U	2.0	1.52	0.76	ug/L	10/31/11	10/31/11
EPA 8260B	1,2-DIBROMOETHANE	0.40 U	1.0	0.40	0.20	ug/L	10/31/11	10/31/11
EPA 8260B	1,2-DICHLOROBENZENE	0.34 U	1.0	0.34	0.17	ug/L	10/31/11	10/31/11
EPA 8260B	1,2-DICHLOROETHANE	0.28 U	1.0	0.28	0.14	ug/L	10/31/11	10/31/11
EPA 8260B	1,2-DICHLOROPROPANE	0.34 U	1.0	0.34	0.17	ug/L	10/31/11	10/31/11
EPA 8260B	1,3-DICHLOROBENZENE	0.22 U	1.0	0.22	0.11	ug/L	10/31/11	10/31/11
EPA 8260B	1,3-DICHLOROPROPENE, TOTAL	0.36 U	1.0	0.36	0.18	ug/L	10/31/11	10/31/11
EPA 8260B	1,4-DICHLOROBENZENE	0.38 U	1.0	0.38	0.19	ug/L	10/31/11	10/31/11
EPA 8260B	2-BUTANONE	1.20 U	10.0	1.20	0.60	ug/L	10/31/11	10/31/11
EPA 8260B	4-METHYL-2-PENTANONE	3.80 U	10.0	3.80	1.90	ug/L	10/31/11	10/31/11
EPA 8260B	ACETONE	1.90 U	10.0	1.90	0.95	ug/L	10/31/11	10/31/11
EPA 8260B	BENZENE	0.32 U	1.0	0.32	0.16	ug/L	10/31/11	10/31/11
EPA 8260B	BROMODICHLOROMETHANE	0.28 U	1.0	0.28	0.14	ug/L	10/31/11	10/31/11
EPA 8260B	BROMOFORM	0.28 U	1.0	0.28	0.14	ug/L	10/31/11	10/31/11
EPA 8260B	BROMOMETHANE	0.48 U	2.0	0.48	0.24	ug/L	10/31/11	10/31/11
EPA 8260B	CARBON TETRACHLORIDE	0.20 U	1.0	0.20	0.10	ug/L	10/31/11	10/31/11
EPA 8260B	CHLOROBENZENE	0.42 U	1.0	0.42	0.21	ug/L	10/31/11	10/31/11
EPA 8260B	CHLORODIBROMOMETHANE	0.38 U	1.0	0.38	0.19	ug/L	10/31/11	10/31/11
EPA 8260B	CHLOROETHANE	0.42 U	1.0	0.42	0.21	ug/L	10/31/11	10/31/11
EPA 8260B	CHLOROFORM	0.14 U	1.0	0.14	0.07	ug/L	10/31/11	10/31/11
EPA 8260B	CHLOROMETHANE	0.84 U	1.0	0.84	0.42	ug/L	10/31/11	10/31/11
EPA 8260B	CIS-1,2-DICHLOROETHENE	0.32 U	1.0	0.32	0.16	ug/L	10/31/11	10/31/11
EPA 8260B	ETHYLBENZENE	0.46 U	1.0	0.46	0.23	ug/L	10/31/11	10/31/11
EPA 8260B	GASOLINE	12.12 U	20.0	12.12	6.06	ug/L	10/31/11	10/31/11
EPA 8260B	HEXACHLOROBUTADIENE	0.38 U	1.0	0.38	0.19	ug/L	10/31/11	10/31/11
EPA 8260B	METHYL TERT-BUTYL ETHER	0.52 U	1.0	0.52	0.26	ug/L	10/31/11	10/31/11
EPA 8260B	METHYLENE CHLORIDE	0.70 U	5.0	0.70	0.35	ug/L	10/31/11	10/31/11

Quant Method: CALLW.M
Run #: 1030C38
Instrument: Chico
Sequence: C111030
Dilution Factor: 1
Initials: ARS

Printed: 12/06/11 3:06:18 PM
APPL-F1-SC-NoMC-REG MDLs

EPA 8260B VOCs + Gas Water

Environet, Inc.
650 Iwilei Rd, #204
Honolulu, HI 96817

APPL Inc.
908 North Temperance Avenue
Clovis, CA 93611

Attn: Stacy Fineran

Project: RED HILL/1022-024

ARF: 66102

Sample ID: ES049

APPL ID: AY49336

Sample Collection Date: 10/24/11

QCG: #86RHB-111030AC-161029

Method	Analyte	Result	LOQ	LOD	DL	Units	Extraction Date	Analysis Date
EPA 8260B	STYRENE	0.50 U	1.0	0.50	0.25	ug/L	10/31/11	10/31/11
EPA 8260B	TETRACHLOROETHENE	0.48 U	1.0	0.48	0.24	ug/L	10/31/11	10/31/11
EPA 8260B	TOLUENE	0.34 U	1.0	0.34	0.17	ug/L	10/31/11	10/31/11
EPA 8260B	TRANS-1,2-DICHLOROETHENE	0.38 U	1.0	0.38	0.19	ug/L	10/31/11	10/31/11
EPA 8260B	TRICHLOROETHENE	0.32 U	1.0	0.32	0.16	ug/L	10/31/11	10/31/11
EPA 8260B	VINYL CHLORIDE	0.46 U	1.0	0.46	0.23	ug/L	10/31/11	10/31/11
EPA 8260B	XYLENES (TOTAL)	0.38 U	1.0	0.38	0.19	ug/L	10/31/11	10/31/11
EPA 8260B	SURROGATE: 1,2-DICHLOROETHANE-	98.1	70-120			%	10/31/11	10/31/11
EPA 8260B	SURROGATE: 4-BROMOFLUOROBEN	102	75-120			%	10/31/11	10/31/11
EPA 8260B	SURROGATE: DIBROMOFLUOROMET	100	85-115			%	10/31/11	10/31/11
EPA 8260B	SURROGATE: TOLUENE-D8 (S)	101	85-120			%	10/31/11	10/31/11

Quant Method: CALLW.M
Run #: 1030C38
Instrument: Chico
Sequence: C111030
Dilution Factor: 1
Initials: ARS

Printed: 12/06/11 3:06:18 PM
APPL-F1-SC-NoMC-REG MDLs

Quantitation Report (QT Reviewed)

Data File : M:\CHICO\DATA\C111030\1030C38W.D Vial: 1
Acq On : 31 Oct 11 15:31 Operator: STC
Sample : AY49336W04 Inst : Chico
Misc : Water 10mLw/ IS&S:10-30/10-26-11 Multiplr: 1.00

Quant Time: Nov 3 12:04 2011 Quant Results File: CALLW.RES

Quant Method : M:\CHICO\DATA\C111030\CALLW.M (RTE Integrator)
Title : METHOD 8260
Last Update : Wed Nov 02 14:33:25 2011
Response via : Initial Calibration
DataAcq Meth : V8260

Internal Standards	R.T.	QIon	Response	Conc Units	Dev(Min)
1) Fluorobenzene (IS)	12.85	96	662144	25.00000 ppb	0.02
55) Chlorobenzene-D5 (IS)	18.05	117	450969	25.00000 ppb	0.00
71) 1,4-Dichlorobenzene-D (IS)	22.26	152	251840	25.00000 ppb	0.02
System Monitoring Compounds					
33) Dibromofluoromethane(S)	11.43	111	443039	25.11762 ppb	0.00
Spiked Amount 25.097			Recovery = 100.083%		
38) 1,2-DCA-D4(S)	12.24	65	373247	23.77157 ppb	0.02
Spiked Amount 24.225			Recovery = 98.129%		
56) Toluene-D8(S)	15.51	98	1662069	26.19352 ppb	0.02
Spiked Amount 25.808			Recovery = 101.494%		
64) 4-Bromofluorobenzene(S)	20.13	95	588080	25.86573 ppb	0.02
Spiked Amount 25.459			Recovery = 101.597%		
Target Compounds					Qvalue
96) Naphthalene	25.95	128	17340	0.56629 ppb	# 89

N.D

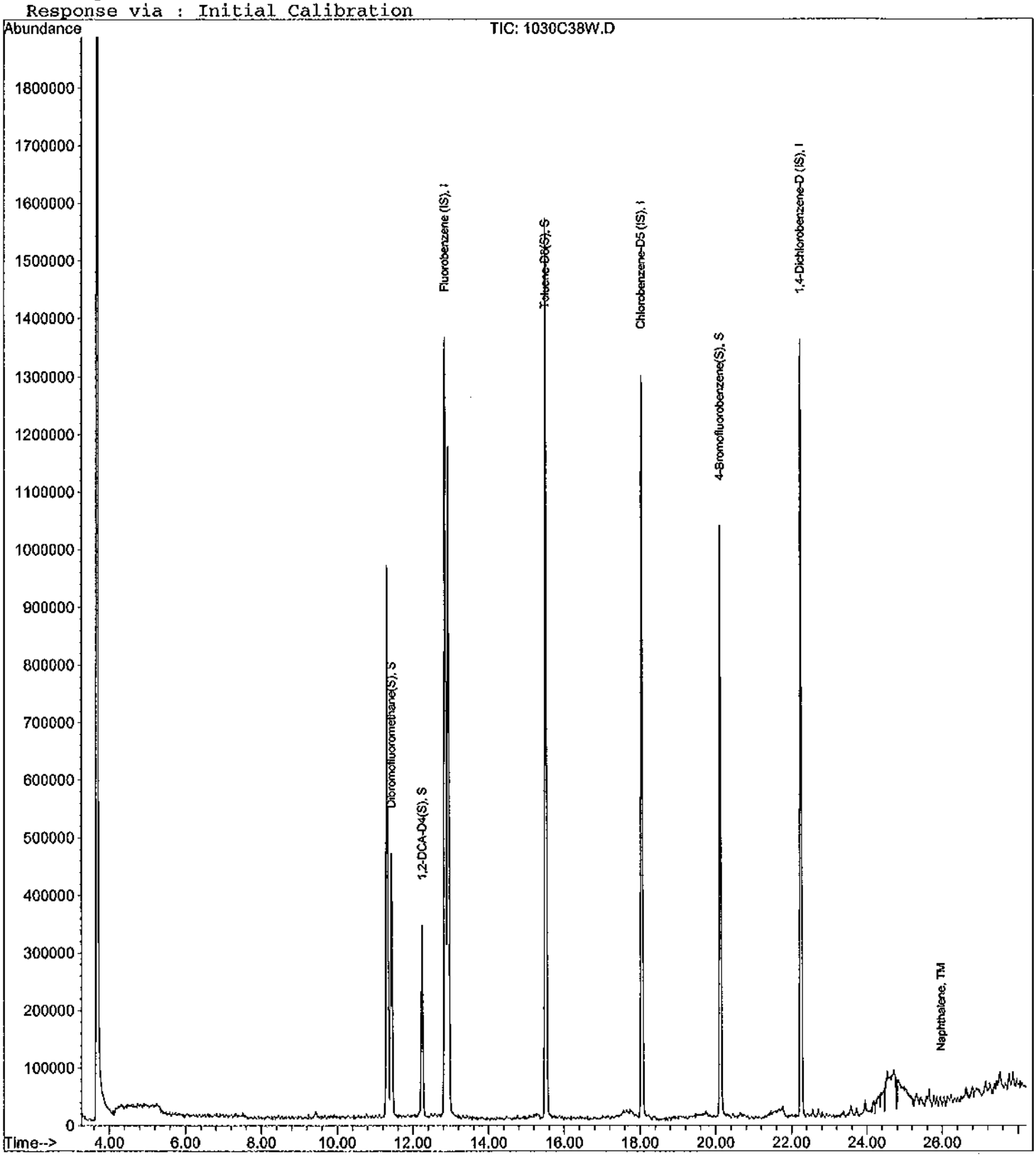
(#) = qualifier out of range (m) = manual integration

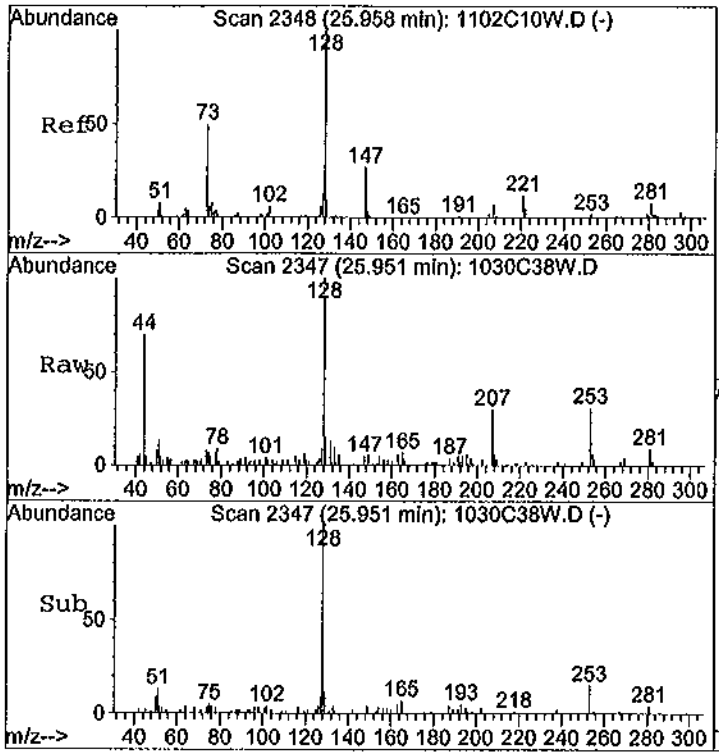
Quantitation Report

Data File : M:\CHICO\DATA\C111030\1030C38W.D Vial: 1
Acq On : 31 Oct 11 15:31 Operator: STC
Sample : AY49336W04 Inst : Chico
Misc : Water 10mLw/ IS&S:10-30/10-26-11 Multiplr: 1.00

Quant Time: Nov 3 12:04 2011 Quant Results File: CALLW.RES

Method : M:\CHICO\DATA\C111030\CALLW.M (RTE Integrator)
Title : METHOD 8260
Last Update : Thu Nov 03 10:27:07 2011
Response via : Initial Calibration

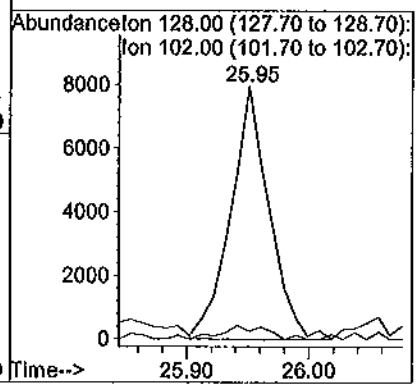




#96
Naphthalene
Concen: 0.56629 ppb
RT: 25.95 min Scan# 2347
Delta R.T. 0.02 min
Lab File: 1030C38W.D
Acq: 31 Oct 11 15:31

Tgt Ion:128 Resp: 17340

Ion	Ratio	Lower	Upper
128	100		
102	3.4	5.1	9.5#



Quantitation Report (QT Reviewed)

Data File : M:\CHICO\DATA\C111030\1030C38W.D Vial: 1
Acq On : 31 Oct 11 15:31 Operator: STC
Sample : AY49336W04 Inst : Chico
Misc : Water 10mLw/ IS&S:10-30/10-26-11 Multiplr: 1.00

Quant Time: Nov 10 10:23 2011 Quant Results File: CGAS.RES

Quant Method : M:\CHICO\DATA\C111030\CGAS.M (RTE Integrator)
Title : METHOD 8260B
Last Update : Thu Nov 03 10:47:02 2011
Response via : Initial Calibration
DataAcq Meth : V8260

Internal Standards	R.T.	QIon	Response	Conc Units	Dev(Min)
1) Fluorobenzene (IS)	12.85	TIC	1349909	25.00000 ppb	0.01
3) Chlorobenzene-D5 (IS)	18.05	TIC	1288292	25.00000 ppb	0.01
4) 1,4-Dichlorobenzene-D (IS)	22.26	TIC	1348071	25.00000 ppb	0.01

System Monitoring Compounds

Target Compounds Qvalue

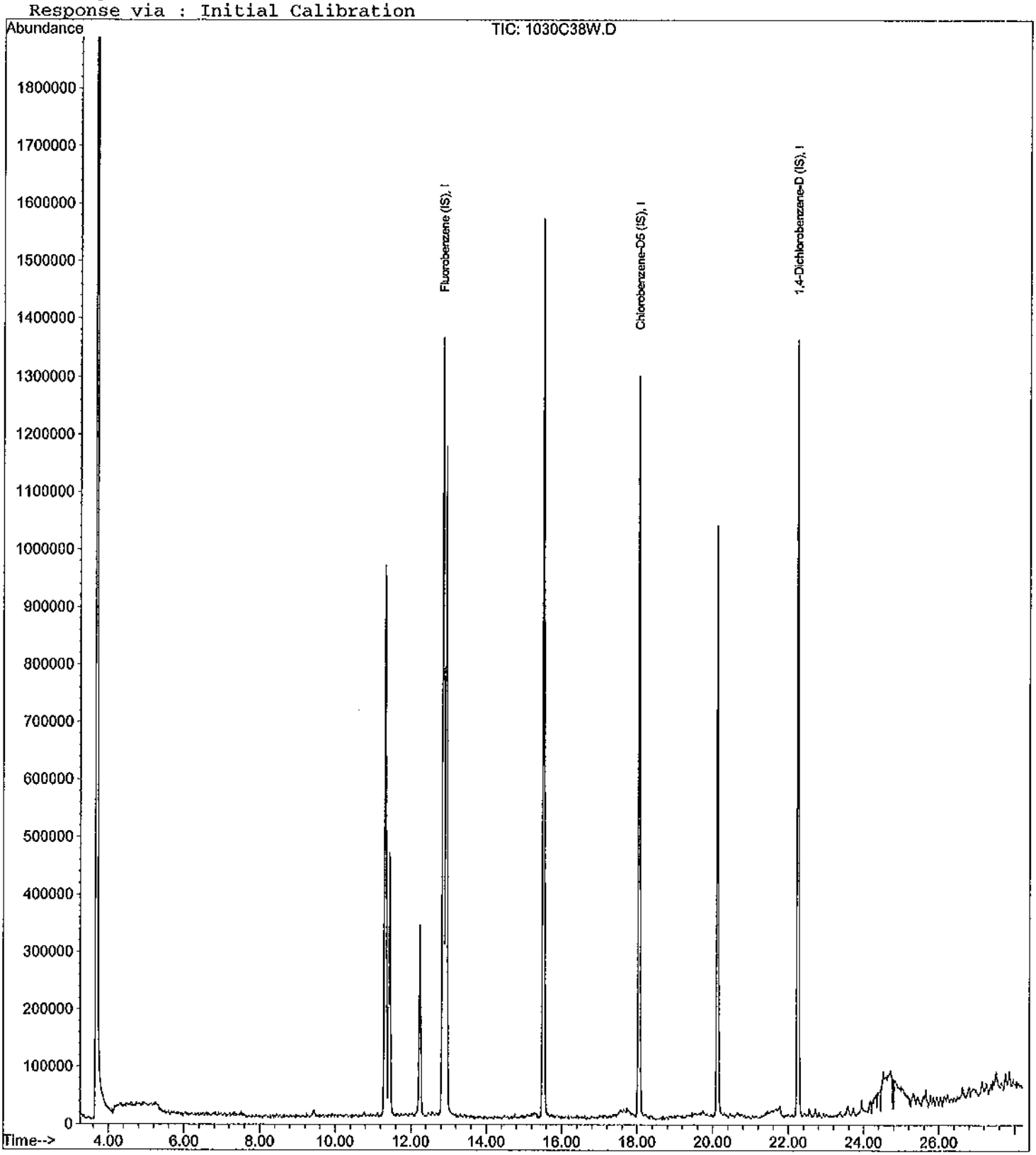
(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : M:\CHICO\DATA\C111030\1030C38W.D Vial: 1
Acq On : 31 Oct 11 15:31 Operator: STC
Sample : AY49336W04 Inst : Chico
Misc : Water 10mLw/ IS&S:10-30/10-26-11 Multiplr: 1.00

Quant Time: Nov 10 10:23 2011 Quant Results File: CGAS.RES

Method : M:\CHICO\DATA\C111030\CGAS.M (RTE Integrator)
Title : METHOD 8260B
Last Update : Thu Nov 03 10:47:02 2011
Response via : Initial Calibration



EPA METHOD 8260B
Volatile Organic Compounds
Calibration Data

APPL, INC.

VOLATILE ORGANIC ANALYSIS BY
EPA METHOD 8260B

Form 6
Initial Calibration

Lab Name: APPL, Inc. SDG No:
Case No: Initial Cal. Date: 10/30/11
Matrix: Instrument: Chico Initials:

			1030C15W.D	1030C16W.D	1030C17W.D	1030C18W.D	1030C19W.D	1030C20W.D	1030C21W.D	1030C22W.D	1030C23W.D					
		Compound	0.3	0.5	1	2	5	10	20	40	100		Avg	%RSD		r
1	I	Fluorobenzene (IS)	ISTD													
2	TM	Dichlorodifluoromethane	0.9991	0.9593	0.9680	0.8471	0.8830	0.9152	0.9504	0.9592	0.8084		0.92	6.8	TM	
3	TM	Freon 114	0.5029	0.5470	0.6167	0.5799	0.5567	0.6260	0.6223	0.6040	0.5364		0.58	7.6	TM	
4	TM**	Chloromethane	1.389	1.171	1.239	1.184	1.061	1.077	1.079	1.076	0.9963		1.1	10	TM**	
5	TM*	Vinyl chloride		0.8227	0.9389	0.7274	0.8810	0.7956	0.7782	0.6322	0.5319		0.76	17	TM*	
6	TML	1,3-Butadiene													TML	
7	TM	Bromomethane	0.5647	0.5370	0.4933	0.6483	0.5131	0.5331	0.5691	0.5732	0.5561		0.55	8.0	TM	
8	TM	Chloroethane		0.7481	0.7782	0.6360	0.6082	0.5955	0.5730	0.5758	0.5302		0.63	14	TM	
9	TM	Dichlorofluoromethane	1.806	1.879	1.897	1.843	1.787	1.725	1.664	1.599	1.493		1.7	7.8	TM	
10	TM	Trichlorofluoromethane	1.219	0.9410	1.059	1.063	1.038	1.037	1.027	1.027	0.9025		1.0	8.5	TM	
11		Acetonitrile	0.0248	0.0305	0.0278	0.0270	0.0288	0.0267	0.0258	0.0269	0.0283		0.03	6.1		
12	TM	Acrolein	0.0160	0.0133	0.0127	0.0117	0.0124	0.0118	0.0112	0.0121	0.0115		0.01	12	TM	
13	TML	Acetone	0.2927	0.4962	0.2742	0.1724	0.1160	0.0970	0.0807	0.0738	0.0705		0.19	77	TML	1.000
14	TML	Freon-113		0.2687	0.6640	0.6219	0.6334	0.6298	0.6085	0.6058	0.5403		0.57	22	TML	0.999
15	TM*	1,1-DCE	0.8684	0.8302	0.7551	0.7075	0.7021	0.6844	0.6445	0.6344	0.5970		0.71	13	TM*	
16	TM	t-Butanol	0.0032	0.0028	0.0031	0.0037	0.0033	0.0035	0.0031	0.0036	0.0043		0.00	13	TM	
17	TML	Methyl Acetate		0.5858	0.4123	0.3223	0.1877	0.2224	0.2002	0.2076	0.2032		0.29	48	TML	1.000
18	TML	Iodomethane			0.1817	0.2265	0.2408	0.2984	0.3981	0.4959	0.6090		0.35	45	TML	0.997
19	TML	Acrylonitrile		0.0466	0.0701	0.0979	0.0824	0.0851	0.0770	0.0761	0.0758		0.08	19	TML	1.000
20	TM	Methylene chloride		0.8304	0.7142	0.7211	0.6635	0.6652	0.6531	0.6094	0.5892		0.68	11	TM	
21	TM	Carbon disulfide	0.8377	0.7169	0.7402	0.7186	0.6675	0.6811	0.6507	0.6311	0.5977		0.69	10	TM	
22	TM	Methyl t-butyl ether (MtBE)	1.160	1.072	1.146	1.130	1.059	1.101	1.037	1.041	0.9630		1.1	5.8	TM	
23	TM	Trans-1,2-DCE	0.9275	0.9146	0.8717	0.8807	0.8240	0.8200	0.7565	0.7483	0.7085		0.83	9.3	TM	
24	TM	Diisopropyl Ether	2.775	2.360	2.465	2.461	2.400	2.425	2.261	2.245	2.073		2.4	8.1	TM	
25	TM**	1,1-DCA	1.448	1.345	1.449	1.466	1.510	1.465	1.402	1.378	1.267		1.4	5.3	TM**	
26	TML	Vinyl Acetate		1.048	0.7027	0.5800	0.5004	0.4421	0.4155	0.4257	0.3841		0.56	40	TML	0.999
27	TM	Ethyl tert Butyl Ether	1.698	1.549	1.596	1.761	1.670	1.751	1.601	1.569	1.460		1.6	6.1	TM	
28	TML	MEK (2-Butanone)		0.5665	0.4588	0.3602	0.2874	0.2964	0.2742	0.2705			0.35	32	TML	1.000
29	TM	Cis-1,2-DCE	1.025	0.9874	0.8510	0.8618	0.8314	0.8432	0.7892	0.7602	0.7089		0.85	12	TM	
30	TM	2,2-Dichloropropane	1.246	1.113	0.9950	1.048	1.020	0.9872	0.9664	0.9254	0.8189		1.0	12	TM	
31	TM*	Chloroform	1.473	1.359	1.352	1.398	1.420	1.399	1.327	1.309	1.208		1.4	5.6	TM*	
32	TM	Bromochloromethane	0.2100	0.2283	0.2743	0.2327	0.2615	0.2451	0.2386	0.2267	0.2148		0.24	8.8	TM	
33	S	Dibromofluoromethane(S)	0.6664	0.6677	0.6754	0.6906	0.6830	0.6822	0.6563	0.6514	0.6207		0.67	3.2	S	
34	TM	1,1,1-TCA	1.307	1.226	1.269	1.241	1.235	1.281	1.263	1.210	1.104		1.2	4.7	TM	
35	TM	Cyclohexane	1.116	1.365	1.215	1.216	1.106	1.144	1.092	1.114	1.000		1.2	8.9	TM	

VOLATILE ORGANIC ANALYSIS BY
EPA METHOD 8260B

Form 6
Initial Calibration

Lab Name: APPL, Inc. SDG No:
Case No: Initial Cal. Date: 10/30/11
Matrix: Instrument: Chico Initials:

		Compound	0.3	0.5	1	2	5	10	20	40	100		Avg	%RSD		r
36	TM	1,1-Dichloropropene	1.089	1.043	1.238	1.061	1.080	1.064	1.039	0.9929	0.9342		1.1	7.8	TM	
37	TML	2,2,4-Trimethylpentane		4.510	2.958	2.198	1.929	1.816	1.761	1.722	1.632		2.3	42	TML	1.000
38	S	1,2-DCA-D4(S)	0.5562	0.6935	0.6364	0.6275	0.6073	0.5850	0.5554	0.5514	0.5227		0.59	9.0	S	
39	TM	Carbon Tetrachloride	0.7745	0.7961	0.7551	0.8397	0.8717	0.9031	0.9401	0.9166	0.8718		0.85	7.6	TM	
40	TM	Tert Amyl Methyl Ether	1.206	1.355	1.238	1.270	1.190	1.245	1.179	1.154	1.119		1.2	5.7	TM	
41	TM	1,2-DCA	0.8312	0.6440	0.6792	0.7198	0.7101	0.7316	0.6883	0.6614	0.6203		0.70	8.8	TM	
42	TM	Benzene	3.278	3.031	3.234	3.139	3.058	3.078	2.926	2.893	2.778		3.0	5.3	TM	
43	TM	TCE	0.7582	0.8575	0.9261	0.9155	0.8745	0.8692	0.8308	0.8139	0.7474		0.84	7.4	TM	
44	TM	2-Pentanone	0.1621	0.1829	0.1696	0.1839	0.1797	0.1849	0.1709	0.1792	0.1751		0.18	4.4	TM	
45	TM*	1,2-Dichloropropane	0.6688	0.7192	0.7502	0.6580	0.7308	0.7195	0.6791	0.6772	0.6288		0.69	5.7	TM*	
46	TM	Bromodichloromethane	0.7189	0.7057	0.8148	0.7719	0.8069	0.8660	0.8196	0.8390	0.7766		0.79	6.7	TM	
47	TM	Methyl Cyclohexane	1.113	1.125	0.9652	0.9330	0.9573	1.003	0.9550	0.9420	0.8802		0.99	8.3	TM	
48	TM	Dibromomethane	0.2413	0.2669	0.2816	0.2846	0.2915	0.3053	0.2877	0.2711	0.2623		0.28	6.8	TM	
49	TM	2-Chloroethyl vinyl ether	0.1532	0.1448	0.1599	0.1939	0.1910	0.1855	0.1751	0.1885	0.1924		0.18	11	TM	
50	TM	1-Bromo-2-chloroethane	0.6355	0.5704	0.5860	0.6008	0.6129	0.6029	0.5729	0.5868	0.5488		0.59	4.3	TM	
51	TM	Cis-1,3-Dichloropropene	0.7822	0.6621	0.7733	0.7420	0.7627	0.7998	0.7723	0.7726	0.7216		0.75	5.5	TM	
52	TM*	Toluene	3.411	3.085	2.935	3.035	3.024	3.066	2.913	2.874	2.698		3.0	6.5	TM*	
53	TM	Trans-1,3-Dichloropropene	0.5191	0.5428	0.4995	0.5430	0.5365	0.5848	0.5511	0.5622	0.5483		0.54	4.5	TM	
54	TM	1,1,2-TCA	0.3181	0.2834	0.2608	0.3010	0.2945	0.3288	0.2891	0.2919	0.2665		0.29	7.5	TM	
55	I	Chlorobenzene-D5 (IS)	ISTD													
56	S	Toluene-D8(S)	3.825	3.742	3.592	3.490	3.642	3.496	3.501	3.233	3.138		3.5	6.3	S	
57	TM	1,2-EDB	0.5186	0.3944	0.4412	0.4259	0.5017	0.5094	0.5156	0.5003	0.4752		0.48	9.4	TM	
58	TM	Tetrachloroethene	1.569	1.513	1.319	1.276	1.308	1.257	1.237	1.102	0.9877		1.3	14	TM	
59	TM	1-Chlorohexane	1.590	1.521	1.436	1.470	1.527	1.492	1.526	1.416	1.343		1.5	5.0	TM	
60	TM	1,1,1,2-Tetrachloroethane	0.6133	0.6996	0.7082	0.7484	0.8608	0.9066	0.9437	0.9002	0.8611		0.80	14	TM	
61	TM	m&p-Xylene	2.262	1.857	1.843	1.755	1.899	1.909	1.977	1.824	1.763		1.9	8.1	TM	
62	TM	o-Xylene	1.911	1.844	1.639	1.684	1.946	1.920	1.973	1.805	1.714		1.8	6.7	TM	
63	TM	Styrene	2.668	2.626	2.667	2.614	2.923	2.903	3.005	2.787	2.612		2.8	5.5	TM	
64	S	4-Bromofluorobenzene(S)	1.386	1.298	1.312	1.224	1.272	1.281	1.253	1.176	1.141		1.3	5.8	S	
65	TM	2-Hexanone		0.2681	0.1979	0.2383	0.2161	0.2363	0.2325	0.2267	0.2142		0.23	9.1	TM	
66	TM	1,3-Dichloropropane	0.9515	0.9439	0.8870	0.9272	0.9571	1.037	0.9690	0.9322	0.8392		0.94	5.8	TM	
67	TM	Dibromochloromethane		0.4957	0.5504	0.5206	0.6049	0.6702	0.6987	0.6794	0.6801		0.61	13	TM	
68	TM**	Chlorobenzene	3.339	2.652	2.608	2.538	2.763	2.780	2.740	2.590	2.437		2.7	9.5	TM**	
69	TM*	Ethylbenzene	5.842	4.934	5.056	4.770	5.229	5.219	5.128	4.773	4.573		5.1	7.3	TM*	
70	TM**L	Bromoform		0.1505	0.1690	0.2072	0.2509	0.2961	0.3266	0.3337	0.3518		0.26	30	TM**L	1.000

VOLATILE ORGANIC ANALYSIS BY
EPA METHOD 8260B

Form 6
Initial Calibration

Lab Name: APPL, Inc. SDG No: _____
Case No: _____ Initial Cal. Date: 10/30/11
Matrix: _____ Instrument: Chico
Initials: _____

		Compound	0.3	0.5	1	2	5	10	20	40	100		Avg	%RSD		r
71	I	1,4-Dichlorobenzene-D (IS)	ISTD													
72	TM	MIBK (methyl isobutyl ketone)		0.7574	0.8768	0.8241	0.6092	0.6987	0.6391	0.6372	0.6432		0.71	14	TM	
73	TM	Isopropylbenzene	9.699	9.039	9.300	9.316	9.115	9.225	9.084	8.340	8.429		9.1	4.7	TM	
74	TM**	1,1,2,2-Tetrachloroethane	0.7160	0.6740	0.7118	0.7852	0.7159	0.8359	0.8186	0.7717	0.7972		0.76	7.4	TM**	
75	TML	1,2,3-Trichloropropane		0.1080	0.1466	0.0826	0.0875	0.1024	0.0817	0.0773	0.0871		0.10	24	TML	0.999
76	TM	t-1,4-Dichloro-2-Butene			0.1384	0.1583	0.1653	0.1826	0.1817	0.1785	0.1992		0.17	12	TM	
77	TM	Bromobenzene	2.191	2.293	2.136	2.173	2.073	2.079	2.023	1.887	1.956		2.1	6.0	TM	
78	TM	n-Propylbenzene	10.7	10.8	11.2	11.8	11.0	11.1	10.9	9.840	10.0		11	5.5	TM	
79	TM	4-Ethyltoluene	8.877	7.431	7.642	7.504	7.338	7.570	7.245	6.748	6.966		7.5	8.0	TM	
80	TM	2-Chlorotoluene	7.018	7.963	7.290	7.456	7.317	7.299	7.188	6.487	6.411		7.2	6.7	TM	
81	TM	1,3,5-Trimethylbenzene	7.727	7.461	7.573	7.610	7.521	7.220	7.414	6.796	6.911		7.4	4.4	TM	
82	TM	4-Chlorotoluene	6.723	5.915	6.587	6.584	6.067	6.161	6.027	5.579	5.836		6.2	6.3	TM	
83	TM	Tert-Butylbenzene	7.797	8.301	7.989	8.311	8.009	8.197	8.168	7.421	7.513		8.0	4.1	TM	
84	TM	1,2,4-Trimethylbenzene	9.214	8.117	7.974	7.683	7.602	7.509	7.430	6.744	6.901		7.7	9.4	TM	
85	TM	Sec-Butylbenzene	9.285	9.224	9.540	9.998	9.615	10.0	10.0	9.119	9.181		9.6	3.9	TM	
86	TM	p-Isopropyltoluene	8.163	8.064	7.913	8.723	8.266	8.400	8.451	7.803	7.874		8.2	3.7	TM	
87	TM	Benzyl Chloride	1.279	1.152	0.9338	1.032	1.032	1.034	1.076	1.092	1.146		1.1	9.0	TM	
88	TM	1,3-DCB	4.734	4.325	4.213	4.479	4.312	4.203	4.279	3.938	3.986		4.3	5.6	TM	
89	TM	1,4-DCB	4.442	3.837	3.983	4.080	3.868	4.040	3.960	3.691	3.802		4.0	5.5	TM	
90	TML	Hexachloroethane		0.4249	0.7118	0.7588	0.9050	1.121	1.321	1.363	1.564		1.0	38	TML	0.999
91	TM	n-Butylbenzene	8.074	7.623	7.277	7.189	6.988	7.038	7.040	6.417	6.597		7.1	7.0	TM	
92	TM	1,2-DCB	3.608	3.335	3.210	3.537	3.348	3.563	3.491	3.215	3.291		3.4	4.5	TM	
93	TML	1,2-Dibromo-3-chloropropane			0.0874	0.0872	0.1047	0.1419	0.1157	0.1281	0.1388		0.11	20	TML	0.999
94	TM	1,2,4-Trichlorobenzene	2.498	2.793	2.514	2.479	2.457	2.481	2.418	2.256	2.277		2.5	6.3	TM	
95	TM	Hexachlorobutadiene		0.5819	0.4282	0.4193	0.3927	0.4607	0.4714	0.4161	0.4105		0.45	13	TM	
96	TM	Naphthalene	3.183	3.265	3.072	3.033	2.834	3.138	3.029	2.875	2.927		3.0	4.7	TM	
97	TM	1,2,3-Trichlorobenzene	1.684	1.977	1.763	2.168	1.821	1.952	1.854	1.766	1.787		1.9	7.9	TM	
98	TML	Dibromochlorobenzene													TML	
99																
100																
101																
102																
103																
104																
105																

Quantitation Report (Not Reviewed)

Data File : M:\CHICO\DATA\C111030\1030C15W.D Vial: 1
Acq On : 30 Oct 11 23:28 Operator: STC
Sample : Voc Std 10-30-11@0.3ug/L Inst : Chico
Misc : Water 10mLw/ IS:10-30-11 Multiplr: 1.00

Quant Time: Dec 2 11:19 2011 Quant Results File: CALLW.RES

Quant Method : M:\CHICO\DATA\C111030\CALLW.M (RTE Integrator)
Title : METHOD 8260
Last Update : Fri Dec 02 11:18:49 2011
Response via : Initial Calibration
DataAcq Meth : V8260

Internal Standards	R.T.	QIon	Response	Conc Units		Dev(Min)
1) Fluorobenzene (IS)	12.83	96	559104	25.00000 ppb		0.00
55) Chlorobenzene-D5 (IS)	18.03	117	374592	25.00000 ppb		-0.01
71) 1,4-Dichlorobenzene-D (IS)	22.24	152	198336	25.00000 ppb		0.00
System Monitoring Compounds						
33) Dibromofluoromethane(S)	11.41	111	8942	0.60039 ppb		-0.01
Spiked Amount 25.097			Recovery	= 2.391%		
38) 1,2-DCA-D4(S)	12.23	65	7464	0.56298 ppb		0.00
Spiked Amount 24.225			Recovery	= 2.324%		
56) Toluene-D8(S)	15.50	98	34391	0.65250 ppb		0.00
Spiked Amount 25.808			Recovery	= 2.526%		
64) 4-Bromofluorobenzene(S)	20.11	95	12464	0.65998 ppb		0.00
Spiked Amount 25.459			Recovery	= 2.592%		
Target Compounds						Qvalue
2) Dichlorodifluoromethane	4.07	85	6703	0.32540 ppb	#	80
3) Freon 114	4.33	85	3374	0.26153 ppb	#	70
4) Chloromethane	4.55	50	9320	0.36511 ppb		94
5) Vinyl chloride	4.83	62	6233	0.36505 ppb	#	76
6) 1,3-Butadiene	4.86	54	426	8.92736 ppb	#	41
7) Bromomethane	5.72	94	3789	0.30570 ppb		70
8) Chloroethane	5.92	64	6243	0.44265 ppb	#	55
9) Dichlorofluoromethane	6.01	67	12120	0.31081 ppb		86
10) Trichlorofluoromethane	6.51	101	8177	0.35336 ppb		78
11) Acetonitrile	7.67	41	8307	13.56274 ug/l		100
12) Acrolein	7.18	56	5382	19.20236 ppb	#	71
13) Acetone	7.26	43	1964	1.22606 ppb	#	75
14) Freon-113	7.48	101	3974	-0.83183 ppb	#	68
15) 1,1-DCE	7.68	96	5826	0.36499 ppb	#	38
16) t-Butanol	7.78	59	1058	13.95861 ppb		96
17) Methyl Acetate	8.17	43	2286	-0.18237 ppb	#	87
18) Iodomethane	8.17	142	479	3.85678 ppb	#	37
19) Acrylonitrile	8.56	53	560	-0.04894 ppb	#	5
20) Methylene chloride	8.46	84	7036	0.46214 ppb	#	60
21) Carbon disulfide	8.57	76	5620	0.36236 ppb		97
22) Methyl t-butyl ether (MtBE	8.91	73	7781	0.32251 ppb	#	64
23) Trans-1,2-DCE	9.09	96	6223	0.38166 ppb		92
24) Diisopropyl Ether	9.75	45	18616	0.34902 ppb	#	73
25) 1,1-DCA	9.78	63	9717	0.30719 ppb		95
26) Vinyl Acetate	9.43	43	9166	-0.47709 ppb	#	81
27) Ethyl tert Butyl Ether	10.45	59	11390	0.31279 ppb		93
28) MEK (2-Butanone)	10.45	43	4667	0.10206 ppb	#	76
29) Cis-1,2-DCE	10.80	96	6876	0.36134 ppb	#	67
30) 2,2-Dichloropropane	10.79	77	8357	0.36882 ppb		95
31) Chloroform	11.08	83	9885	0.32484 ppb		86
32) Bromochloromethane	11.31	128	1409	0.26596 ppb	#	1
34) 1,1,1-TCA	11.83	97	8769	0.31687 ppb	#	72
35) Cyclohexane	11.98	56	7488	0.29062 ppb	#	74
36) 1,1-Dichloropropene	12.10	75	7308	0.30819 ppb		86
37) 2,2,4-Trimethylpentane	12.17	57	48085	0.17131 ppb	#	87
39) Carbon Tetrachloride	12.31	117	5196	0.27267 ppb	#	87

(#) = qualifier out of range (m) = manual integration

Quantitation Report (Not Reviewed)

Data File : M:\CHICO\DATA\C111030\1030C15W.D Vial: 1
Acq On : 30 Oct 11 23:28 Operator: STC
Sample : Voc Std 10-30-11@0.3ug/L Inst : Chico
Misc : Water 10mLw/ IS:10-30-11 Multiplr: 1.00

Quant Time: Dec 2 11:19 2011 Quant Results File: CALLW.RES

Quant Method : M:\CHICO\DATA\C111030\CALLW.M (RTE Integrator)
Title : METHOD 8260
Last Update : Fri Dec 02 11:18:49 2011
Response via : Initial Calibration
DataAcq Meth : V8260

	Compound	R.T.	QIon	Response	Conc Unit		Qvalue
40)	Tert Amyl Methyl Ether	12.35	73	8093	0.29724 ppb	#	82
41)	1,2-DCA	12.36	62	5577	0.35705 ppb	#	79
42)	Benzene	12.50	78	21995	0.32285 ppb		99
43)	TCE	13.53	95	5087	0.26961 ppb		88
44)	2-Pentanone	13.19	43	54377	13.77755 ppb		91
45)	1,2-Dichloropropane	13.76	63	4487	0.28977 ppb	#	81
46)	Bromodichloromethane	14.11	83	4823	0.27263 ppb	#	84
47)	Methyl Cyclohexane	13.81	83	7465	0.33857 ppb		81
48)	Dibromomethane	14.15	93	1619	0.26143 ppb	#	56
49)	2-Chloroethyl vinyl ether	14.57	63	1028	0.26114 ppb	#	71
50)	1-Bromo-2-chloroethane	14.89	63	4264	0.32272 ppb	#	76
51)	Cis-1,3-Dichloropropene	15.00	75	5248	0.31110 ppb		96
52)	Toluene	15.64	91	22887	0.34061 ppb		88
53)	Trans-1,3-Dichloropropene	15.80	75	3483	0.28679 ppb	#	72
54)	1,1,2-TCA	16.08	83	2134	0.32602 ppb		86
57)	1,2-EDB	17.33	107	2331	0.32696 ppb	#	72
58)	Tetrachloroethene	16.79	164	7051	0.36612 ppb	#	74
59)	1-Chlorohexane	17.71	91	7147	0.32226 ppb	#	70
60)	1,1,1,2-Tetrachloroethane	18.16	131	2757	0.22867 ppb		83
61)	m&p-Xylene	18.35	106	20335	0.71474 ppb		85
62)	o-Xylene	19.11	106	8588	0.31386 ppb		68
63)	Styrene	19.12	104	11993	0.29039 ppb		98
65)	2-Hexanone	16.12	43	2535	0.73958 ppb	#	73
66)	1,3-Dichloropropane	16.49	76	4277	0.30423 ppb		96
67)	Dibromochloromethane	16.98	129	1907	0.20779 ppb	#	39
68)	Chlorobenzene	18.10	112	15008	0.36875 ppb	#	68
69)	Ethylbenzene	18.21	91	26259	0.34648 ppb		89
70)	Bromoform	19.63	173	532	1.23415 ppb	#	37
72)	MIBK (methyl isobutyl keto	14.68	43	7070	1.25392 ppb		81
73)	Isopropylbenzene	19.72	105	23083	0.32111 ppb	#	81
74)	1,1,2,2-Tetrachloroethane	19.88	83	1704	0.28318 ppb	#	79
75)	1,2,3-Trichloropropane	20.14	110	346	0.59881 ppb	#	25
76)	t-1,4-Dichloro-2-Butene	20.24	53	170	0.12459 ppb	#	62
77)	Bromobenzene	20.49	156	5215	0.31451 ppb	#	54
78)	n-Propylbenzene	20.44	91	25363	0.29576 ppb		97
79)	4-Ethyltoluene	20.63	105	21128	0.35603 ppb		93
80)	2-Chlorotoluene	20.74	91	16704	0.29411 ppb		87
81)	1,3,5-Trimethylbenzene	20.72	105	18391	0.31501 ppb		98
82)	4-Chlorotoluene	20.81	91	16002	0.32720 ppb	#	74
83)	Tert-Butylbenzene	21.36	119	18557	0.29358 ppb		92
84)	1,2,4-Trimethylbenzene	21.41	105	21930	0.35965 ppb		91
85)	Sec-Butylbenzene	21.75	105	22099	0.29154 ppb		86
86)	p-Isopropyltoluene	21.99	119	19428	0.29923 ppb		97
87)	Benzyl Chloride	22.41	91	3043	0.35307 ppb	#	69
88)	1,3-DCB	22.12	146	11266	0.33224 ppb		92
89)	1,4-DCB	22.29	146	10573	0.33594 ppb		95
90)	Hexachloroethane	23.59	117	857	2.03197 ppb	#	50
91)	n-Butylbenzene	22.70	91	19217	0.33934 ppb		91
92)	1,2-DCB	22.92	146	8586	0.31833 ppb	#	78
93)	1,2-Dibromo-3-chloropropan	24.14	155	141	1.41683 ppb	#	27
94)	1,2,4-Trichlorobenzene	25.58	180	5945	0.30417 ppb	#	86

(#) = qualifier out of range (m) = manual integration

Quantitation Report (Not Reviewed)

Data File : M:\CHICO\DATA\C111030\1030C15W.D Vial: 1
Acq On : 30 Oct 11 23:28 Operator: STC
Sample : Voc Std 10-30-11@0.3ug/L Inst : Chico
Misc : Water 10mLw/ IS:10-30-11 Multiplr: 1.00

Quant Time: Dec 2 11:19 2011 Quant Results File: CALLW.RES

Quant Method : M:\CHICO\DATA\C111030\CALLW.M (RTE Integrator)
Title : METHOD 8260
Last Update : Fri Dec 02 11:18:49 2011
Response via : Initial Calibration
DataAcq Meth : V8260

Compound	R.T.	QIon	Response	Conc	Unit		Qvalue
95) Hexachlorobutadiene	25.83	223	1530	0.43087	ppb	#	68
96) Naphthalene	25.92	128	7576	0.31416	ppb	#	79
97) 1,2,3-Trichlorobenzene	26.29	180	4008	0.27110	ppb		85

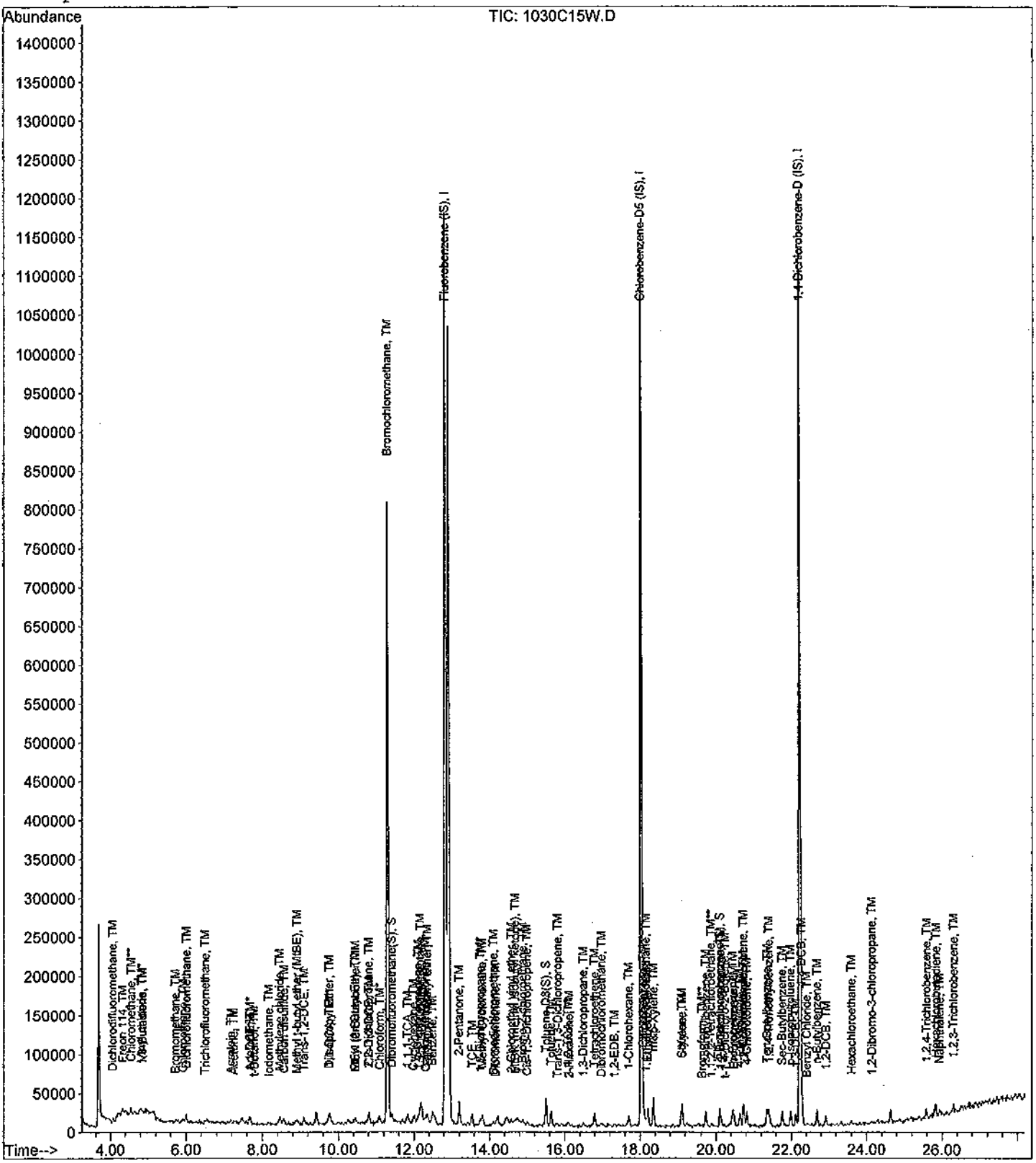
(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : M:\CHICO\DATA\C111030\1030C15W.D Vial: 1
Acq On : 30 Oct 11 23:28 Operator: STC
Sample : Voc Std 10-30-11@0.3ug/L Inst : Chico
Misc : Water 10mLw/ IS:10-30-11 Multiplr: 1.00

Quant Time: Dec 2 11:19 2011 Quant Results File: CALLW.RES

Method : M:\CHICO\DATA\C111030\CALLW.M (RTE Integrator)
Title : METHOD 8260
Last Update : Fri Dec 02 11:18:49 2011
Response via : Initial Calibration



Quantitation Report (Not Reviewed)

Data File : M:\CHICO\DATA\C111030\1030C16W.D Vial: 1
Acq On : 31 Oct 11 00:11 Operator: STC
Sample : Voc Std 10-30-11@0.5ug/L Inst : Chico
Misc : Water 10mLw/ IS:10-30-11 Multiplr: 1.00

Quant Time: Dec 2 11:19 2011 Quant Results File: CALLW.RES

Quant Method : M:\CHICO\DATA\C111030\CALLW.M (RTE Integrator)
Title : METHOD 8260
Last Update : Fri Dec 02 11:18:49 2011
Response via : Initial Calibration
DataAcq Meth : V8260

Internal Standards	R.T.	QIon	Response	Conc Units		Dev(Min)
1) Fluorobenzene (IS)	12.84	96	564160	25.00000 ppb		0.00
55) Chlorobenzene-D5 (IS)	18.03	117	384000	25.00000 ppb		0.00
71) 1,4-Dichlorobenzene-D (IS)	22.24	152	199104	25.00000 ppb		0.00
System Monitoring Compounds						
33) Dibromofluoromethane(S)	11.43	111	15067	1.00257 ppb		0.00
Spiked Amount 25.097			Recovery =	3.996%		
38) 1,2-DCA-D4(S)	12.23	65	15649	1.16976 ppb		0.00
Spiked Amount 24.225			Recovery =	4.830%		
56) Toluene-D8(S)	15.51	98	57480	1.06384 ppb		0.01
Spiked Amount 25.808			Recovery =	4.123%		
64) 4-Bromofluorobenzene(S)	20.10	95	19938	1.02988 ppb		0.00
Spiked Amount 25.459			Recovery =	4.046%		
Target Compounds						Qvalue
2) Dichlorodifluoromethane	4.07	85	10824	0.52075 ppb		95
3) Freon 114	4.33	85	6172	0.47412 ppb		98
4) Chloromethane	4.55	50	13218	0.51317 ppb	#	72
5) Vinyl chloride	4.81	62	9283	0.53880 ppb	#	69
6) 1,3-Butadiene	4.79	54	547	11.36033 ppb		97
7) Bromomethane	5.73	94	6059	0.48447 ppb		87
8) Chloroethane	5.91	64	8441	0.59313 ppb	#	74
9) Dichlorofluoromethane	6.02	67	21197	0.53871 ppb	#	80
10) Trichlorofluoromethane	6.51	101	10617	0.45468 ppb		86
11) Acetonitrile	7.65	41	17187	27.80953 ug/l		100
12) Acrolein	7.15	56	7488	26.47690 ppb		95
13) Acetone	7.29	43	5599	3.46393 ppb		93
14) Freon-113	7.43	101	3032	-0.91166 ppb	#	70
15) 1,1-DCE	7.68	96	9367	0.58157 ppb		86
16) t-Butanol	7.75	59	1553	20.30572 ppb		100
17) Methyl Acetate	8.18	43	6610	0.76232 ppb	#	81
18) Iodomethane	8.16	142	1888	3.95744 ppb	#	59
19) Acrylonitrile	8.56	53	526	-0.07185 ppb	#	71
20) Methylene chloride	8.48	84	9369	0.60986 ppb		86
21) Carbon disulfide	8.56	76	8089	0.51688 ppb		94
22) Methyl t-butyl ether (MtBE	8.89	73	12097	0.49691 ppb		93
23) Trans-1,2-DCE	9.08	96	10320	0.62725 ppb		75
24) Diisopropyl Ether	9.74	45	26624	0.49469 ppb	#	77
25) 1,1-DCA	9.79	63	15171	0.47532 ppb	#	86
26) Vinyl Acetate	9.43	43	11823	-0.17906 ppb	#	82
27) Ethyl tert Butyl Ether	10.44	59	17474	0.47557 ppb		99
28) MEK (2-Butanone)	10.44	43	6392	0.38220 ppb	#	76
29) Cis-1,2-DCE	10.81	96	11141	0.58022 ppb	#	62
30) 2,2-Dichloropropane	10.83	77	12554	0.54909 ppb		96
31) Chloroform	11.09	83	15337	0.49949 ppb		92
32) Bromochloromethane	11.30	128	2576	0.48189 ppb	#	33
34) 1,1,1-TCA	11.83	97	13833	0.49538 ppb	#	82
35) Cyclohexane	11.99	56	15403	0.59245 ppb		78
36) 1,1-Dichloropropene	12.10	75	11774	0.49209 ppb	#	84
37) 2,2,4-Trimethylpentane	12.18	57	50886	0.23590 ppb		93
39) Carbon Tetrachloride	12.29	117	8982	0.46713 ppb		86

(#) = qualifier out of range (m) = manual integration

Quantitation Report (Not Reviewed)

Data File : M:\CHICO\DATA\C111030\1030C16W.D Vial: 1
Acq On : 31 Oct 11 00:11 Operator: STC
Sample : Voc Std 10-30-11@0.5ug/L Inst : Chico
Misc : Water 10mLw/ IS:10-30-11 Multiplr: 1.00

Quant Time: Dec 2 11:19 2011 Quant Results File: CALLW.RES

Quant Method : M:\CHICO\DATA\C111030\CALLW.M (RTE Integrator)
Title : METHOD 8260
Last Update : Fri Dec 02 11:18:49 2011
Response via : Initial Calibration
DataAcq Meth : V8260

Compound	R.T.	QIon	Response	Conc Unit		Qvalue
40) Tert Amyl Methyl Ether	12.34	73	15291	0.55657 ppb	#	82
41) 1,2-DCA	12.36	62	7266	0.46101 ppb		99
42) Benzene	12.50	78	34203	0.49755 ppb	#	86
43) TCE	13.53	95	9675	0.50818 ppb	#	76
44) 2-Pentanone	13.21	43	103205	25.91480 ppb		91
45) 1,2-Dichloropropane	13.77	63	8115	0.51936 ppb		98
46) Bromodichloromethane	14.11	83	7962	0.44603 ppb	#	98
47) Methyl Cyclohexane	13.81	83	12697	0.57069 ppb		99
48) Dibromomethane	14.16	93	3011	0.48185 ppb	#	68
49) 2-Chloroethyl vinyl ether	14.58	63	1634	0.41135 ppb	#	59
50) 1-Bromo-2-chloroethane	14.88	63	6436	0.48274 ppb	#	66
51) Cis-1,3-Dichloropropene	15.00	75	7471	0.43891 ppb	#	75
52) Toluene	15.63	91	34808	0.51338 ppb		87
53) Trans-1,3-Dichloropropene	15.80	75	6124	0.49974 ppb	#	83
54) 1,1,2-TCA	16.08	83	3198	0.48419 ppb	#	78
57) 1,2-EDB	17.33	107	3029	0.41446 ppb	#	80
58) Tetrachloroethene	16.78	164	11617	0.58843 ppb		81
59) 1-Chlorohexane	17.70	91	11679	0.51371 ppb		93
60) 1,1,1,2-Tetrachloroethane	18.16	131	5373	0.43472 ppb		80
61) m&p-Xylene	18.36	106	28523	0.97798 ppb		92
62) o-Xylene	19.10	106	14161	0.50486 ppb		88
63) Styrene	19.13	104	20171	0.47643 ppb		89
65) 2-Hexanone	16.11	43	2059	0.58599 ppb	#	78
66) 1,3-Dichloropropane	16.50	76	7249	0.50300 ppb		84
67) Dibromochloromethane	16.95	129	3807	0.40465 ppb		74
68) Chlorobenzene	18.11	112	20370	0.48823 ppb	#	81
69) Ethylbenzene	18.21	91	37896	0.48777 ppb		98
70) Bromoform	19.62	173	1156	1.34654 ppb	#	37
72) MIBK (methyl isobutyl keto	14.67	43	3016	0.53285 ppb	#	51
73) Isopropylbenzene	19.73	105	35995	0.49881 ppb		90
74) 1,1,2,2-Tetrachloroethane	19.89	83	2684	0.44432 ppb	#	86
75) 1,2,3-Trichloropropane	20.15	110	430	0.71972 ppb	#	57
76) t-1,4-Dichloro-2-Butene	20.21	53	1066	0.77827 ppb	#	42
77) Bromobenzene	20.47	156	9129	0.54844 ppb	#	86
78) n-Propylbenzene	20.44	91	42985	0.49932 ppb		89
79) 4-Ethyltoluene	20.63	105	29592	0.49673 ppb		84
80) 2-Chlorotoluene	20.73	91	31711	0.55619 ppb	#	79
81) 1,3,5-Trimethylbenzene	20.71	105	29710	0.50692 ppb		88
82) 4-Chlorotoluene	20.82	91	23555	0.47979 ppb		86
83) Tert-Butylbenzene	21.36	119	33054	0.52092 ppb		95
84) 1,2,4-Trimethylbenzene	21.41	105	32321	0.52802 ppb	#	68
85) Sec-Butylbenzene	21.76	105	36729	0.48267 ppb		87
86) p-Isopropyltoluene	22.00	119	32111	0.49266 ppb		93
87) Benzyl Chloride	22.43	91	4589	0.53039 ppb		91
88) 1,3-DCB	22.11	146	17221	0.50589 ppb	#	84
89) 1,4-DCB	22.29	146	15280	0.48363 ppb		95
90) Hexachloroethane	23.59	117	1692	2.09850 ppb	#	60
91) n-Butylbenzene	22.70	91	30355	0.53395 ppb		90
92) 1,2-DCB	22.92	146	13282	0.49053 ppb		92
93) 1,2-Dibromo-3-chloropropan	24.16	155	468	1.71132 ppb	#	1
94) 1,2,4-Trichlorobenzene	25.58	180	11120	0.56675 ppb		82

(#) = qualifier out of range (m) = manual integration

Quantitation Report (Not Reviewed)

Data File : M:\CHICO\DATA\C111030\1030C16W.D Vial: 1
Acq On : 31 Oct 11 00:11 Operator: STC
Sample : Voc Std 10-30-11@0.5ug/L Inst : Chico
Misc : Water 10mLw/ IS:10-30-11 Multiplr: 1.00

Quant Time: Dec 2 11:19 2011 Quant Results File: CALLW.RES

Quant Method : M:\CHICO\DATA\C111030\CALLW.M (RTE Integrator)
Title : METHOD 8260
Last Update : Fri Dec 02 11:18:49 2011
Response via : Initial Calibration
DataAcq Meth : V8260

Compound	R.T.	QIon	Response	Conc	Unit		Qvalue
95) Hexachlorobutadiene	25.83	223	2317	0.64998	ppb	#	59
96) Naphthalene	25.93	128	13003	0.53713	ppb	#	92
97) 1,2,3-Trichlorobenzene	26.30	180	7873	0.53048	ppb	#	70

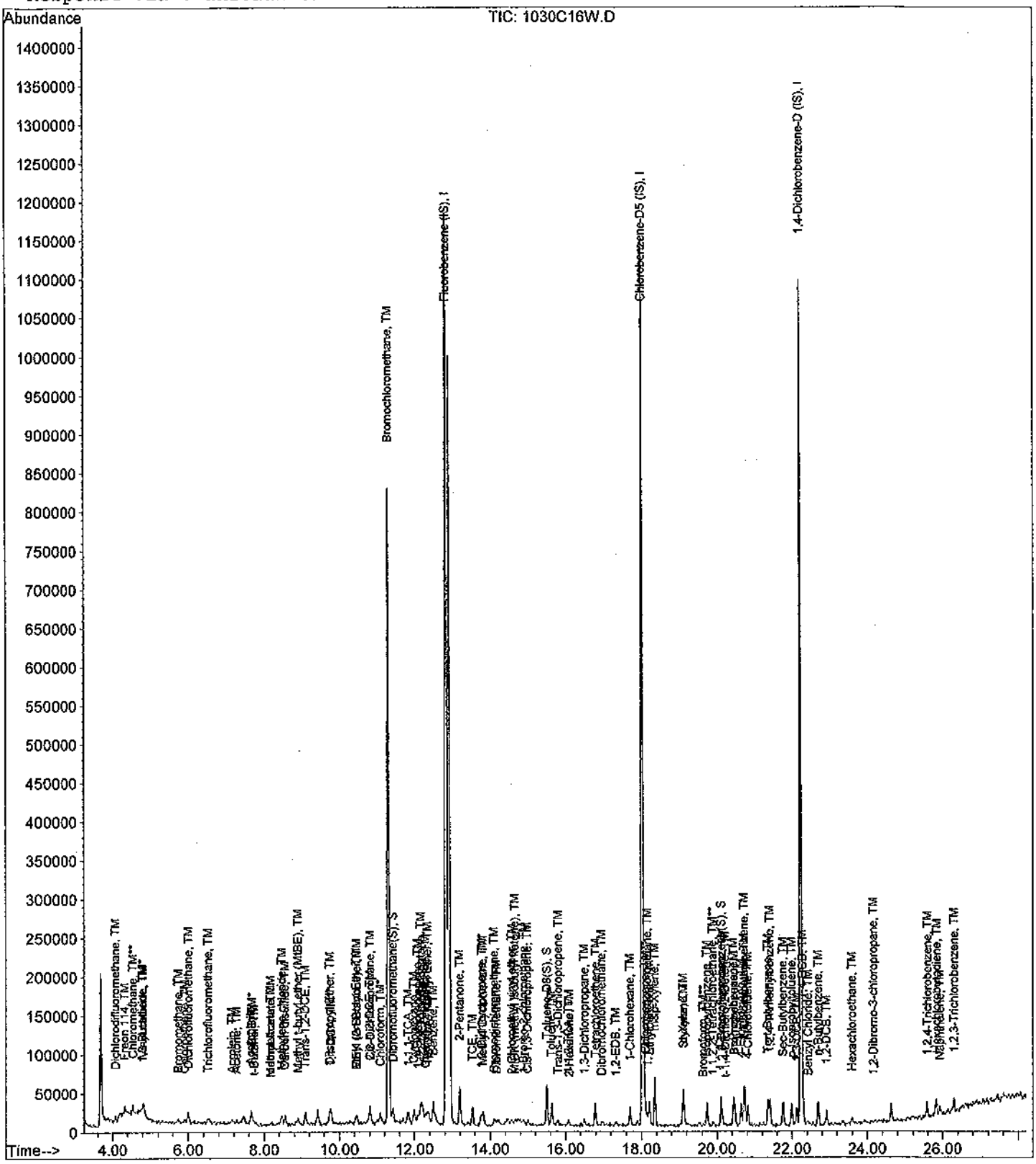
(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : M:\CHICO\DATA\C111030\1030C16W.D Vial: 1
Acq On : 31 Oct 11 00:11 Operator: STC
Sample : Voc Std 10-30-11@0.5ug/L Inst : Chico
Misc : Water 10mLw/ IS:10-30-11 Multiplr: 1.00

Quant Time: Dec 2 11:19 2011 Quant Results File: CALLW.RES

Method : M:\CHICO\DATA\C111030\CALLW.M (RTE Integrator)
Title : METHOD 8260
Last Update : Fri Dec 02 11:18:49 2011
Response via : Initial Calibration



Quantitation Report (Not Reviewed)

Data File : M:\CHICO\DATA\C111030\1030C17W.D Vial: 1
Acq On : 31 Oct 11 00:54 Operator: STC
Sample : Voc Std 10-30-11@1.0ug/L Inst : Chico
Misc : Water 10mLw/ IS:10-30-11 Multiplr: 1.00

Quant Time: Dec 2 11:19 2011 Quant Results File: CALLW.RES

Quant Method : M:\CHICO\DATA\C111030\CALLW.M (RTE Integrator)
Title : METHOD 8260
Last Update : Fri Dec 02 11:18:49 2011
Response via : Initial Calibration
DataAcq Meth : V8260

Internal Standards	R.T.	QIon	Response	Conc	Units		Dev(Min)
1) Fluorobenzene (IS)	12.84	96	539200	25.00000	ppb		0.00
55) Chlorobenzene-D5 (IS)	18.03	117	383872	25.00000	ppb		0.00
71) 1,4-Dichlorobenzene-D (IS)	22.24	152	197760	25.00000	ppb		0.00
System Monitoring Compounds							
33) Dibromofluoromethane(S)	11.42	111	29134	2.02833	ppb		0.00
Spiked Amount 25.097			Recovery = 8.081%				
38) 1,2-DCA-D4(S)	12.22	65	27452	2.14703	ppb		0.00
Spiked Amount 24.225			Recovery = 8.863%				
56) Toluene-D8(S)	15.50	98	110307	2.04225	ppb		0.00
Spiked Amount 25.808			Recovery = 7.912%				
64) 4-Bromofluorobenzene(S)	20.11	95	40278	2.08121	ppb		0.00
Spiked Amount 25.459			Recovery = 8.174%				
Target Compounds							Qvalue
2) Dichlorodifluoromethane	4.07	85	20878	1.05095	ppb	#	77
3) Freon 114	4.34	85	13301	1.06904	ppb		79
4) Chloromethane	4.55	50	26727	1.08568	ppb		96
5) Vinyl chloride	4.82	62	20250	1.22976	ppb		92
6) 1,3-Butadiene	4.82	54	298	6.47548	ppb	#	63
7) Bromomethane	5.73	94	10639	0.89005	ppb		88
8) Chloroethane	5.91	64	16785	1.23404	ppb		95
9) Dichlorofluoromethane	6.01	67	40917	1.08801	ppb		99
10) Trichlorofluoromethane	6.54	101	22833	1.02311	ppb	#	73
11) Acetonitrile	7.65	41	29961	50.72269	ug/l		100
12) Acrolein	7.16	56	13727	50.78428	ppb		87
13) Acetone	7.28	43	5913	3.82754	ppb	#	80
14) Freon-113	7.47	101	14321	0.06398	ppb	#	85
15) 1,1-DCE	7.68	96	16287	1.05802	ppb	#	61
16) t-Butanol	7.76	59	3312	45.30953	ppb		93
17) Methyl Acetate	8.19	43	8892	1.35363	ppb		91
18) Iodomethane	8.17	142	3919	4.11599	ppb	#	77
19) Acrylonitrile	8.55	53	1511	0.54689	ppb	#	42
20) Methylene chloride	8.47	84	15404	1.04912	ppb	#	71
21) Carbon disulfide	8.56	76	15964	1.06731	ppb		97
22) Methyl t-butyl ether (MtBE	8.89	73	24721	1.06247	ppb	#	90
23) Trans-1,2-DCE	9.10	96	18800	1.19557	ppb	#	91
24) Diisopropyl Ether	9.75	45	53157	1.03341	ppb		96
25) 1,1-DCA	9.78	63	31252	1.02448	ppb	#	91
26) Vinyl Acetate	9.41	43	15155	0.28793	ppb	#	78
27) Ethyl tert Butyl Ether	10.44	59	34422	0.98019	ppb		99
28) MEK (2-Butanone)	10.44	43	9896	1.04178	ppb	#	90
29) Cis-1,2-DCE	10.81	96	18355	1.00018	ppb		79
30) 2,2-Dichloropropane	10.81	77	21461	0.98211	ppb		96
31) Chloroform	11.09	83	29160	0.99364	ppb		95
32) Bromochloromethane	11.32	128	5916	1.15793	ppb		90
34) 1,1,1-TCA	11.83	97	27369	1.02550	ppb		88
35) Cyclohexane	12.00	56	26209	1.05474	ppb		95
36) 1,1-Dichloropropene	12.09	75	26709	1.16796	ppb	#	85
37) 2,2,4-Trimethylpentane	12.17	57	63801	0.66906	ppb		93
39) Carbon Tetrachloride	12.29	117	16287	0.88625	ppb		97

(#) = qualifier out of range (m) = manual integration

Quantitation Report (Not Reviewed)

Data File : M:\CHICO\DATA\C111030\1030C17W.D Vial: 1
Acq On : 31 Oct 11 00:54 Operator: STC
Sample : Voc Std 10-30-11@1.0ug/L Inst : Chico
Misc : Water 10mLw/ IS:10-30-11 Multiplr: 1.00

Quant Time: Dec 2 11:19 2011 Quant Results File: CALLW.RES

Quant Method : M:\CHICO\DATA\C111030\CALLW.M (RTE Integrator)
Title : METHOD 8260
Last Update : Fri Dec 02 11:18:49 2011
Response via : Initial Calibration
DataAcq Meth : V8260

	Compound	R.T.	QIon	Response	Conc	Unit		Qvalue
40)	Tert Amyl Methyl Ether	12.35	73	26700	1.01682	ppb		98
41)	1,2-DCA	12.38	62	14649	0.97246	ppb	#	86
42)	Benzene	12.49	78	69760	1.06177	ppb	#	93
43)	TCE	13.53	95	19974	1.09770	ppb		87
44)	2-Pentanone	13.20	43	182845	48.03774	ppb		99
45)	1,2-Dichloropropane	13.77	63	16180	1.08346	ppb		97
46)	Bromodichloromethane	14.12	83	17574	1.03007	ppb	#	98
47)	Methyl Cyclohexane	13.82	83	20818	0.97903	ppb		89
48)	Dibromomethane	14.16	93	6074	1.01701	ppb	#	71
49)	2-Chloroethyl vinyl ether	14.58	63	3448	0.90821	ppb		94
50)	1-Bromo-2-chloroethane	14.88	63	12639	0.99189	ppb	#	62
51)	Cis-1,3-Dichloropropene	15.01	75	16679	1.02522	ppb		86
52)	Toluene	15.63	91	63296	0.97675	ppb		83
53)	Trans-1,3-Dichloropropene	15.81	75	10774	0.91989	ppb		92
54)	1,1,2-TCA	16.09	83	5625	0.89108	ppb		89
57)	1,2-EDB	17.32	107	6774	0.92719	ppb	#	92
58)	Tetrachloroethene	16.79	164	20256	1.02636	ppb		96
59)	1-Chlorohexane	17.70	91	22047	0.97009	ppb		95
60)	1,1,1,2-Tetrachloroethane	18.16	131	10874	0.88009	ppb		87
61)	m&p-Xylene	18.35	106	56585	1.94079	ppb		87
62)	o-Xylene	19.10	106	25167	0.89753	ppb		93
63)	Styrene	19.12	104	40956	0.96769	ppb		88
65)	2-Hexanone	16.10	43	3038	0.86490	ppb	#	70
66)	1,3-Dichloropropane	16.49	76	13620	0.94539	ppb		85
67)	Dibromochloromethane	16.98	129	8452	0.89867	ppb	#	59
68)	Chlorobenzene	18.10	112	40045	0.96012	ppb		92
69)	Ethylbenzene	18.22	91	77629	0.99952	ppb		96
70)	Bromoform	19.64	173	2595	1.61154	ppb	#	81
72)	MIBK (methyl isobutyl keto	14.66	43	6936	1.23373	ppb	#	66
73)	Isopropylbenzene	19.74	105	73566	1.02638	ppb	#	82
74)	1,1,2,2-Tetrachloroethane	19.90	83	5631	0.93852	ppb		82
75)	1,2,3-Trichloropropane	20.15	110	1160	1.79904	ppb	#	72
76)	t-1,4-Dichloro-2-Butene	20.24	53	1095	0.80488	ppb	#	13
77)	Bromobenzene	20.48	156	16894	1.02183	ppb		83
78)	n-Propylbenzene	20.44	91	88461	1.03456	ppb		99
79)	4-Ethyltoluene	20.64	105	60450	1.02161	ppb		93
80)	2-Chlorotoluene	20.73	91	57667	1.01832	ppb		95
81)	1,3,5-Trimethylbenzene	20.71	105	59903	1.02902	ppb		95
82)	4-Chlorotoluene	20.81	91	52108	1.06859	ppb		96
83)	Tert-Butylbenzene	21.35	119	63199	1.00276	ppb		94
84)	1,2,4-Trimethylbenzene	21.42	105	63077	1.03747	ppb		94
85)	Sec-Butylbenzene	21.76	105	75467	0.99848	ppb		94
86)	p-Isopropyltoluene	21.99	119	62593	0.96686	ppb		95
87)	Benzyl Chloride	22.43	91	7387	0.85958	ppb	#	79
88)	1,3-DCB	22.14	146	33327	0.98569	ppb		96
89)	1,4-DCB	22.30	146	31510	1.00410	ppb		98
90)	Hexachloroethane	23.60	117	5631	2.41666	ppb	#	75
91)	n-Butylbenzene	22.70	91	57562	1.01941	ppb		93
92)	1,2-DCB	22.92	146	25391	0.94412	ppb	#	84
93)	1,2-Dibromo-3-chloropropan	24.15	155	691	1.91673	ppb	#	73
94)	1,2,4-Trichlorobenzene	25.59	180	19884	1.02031	ppb		78

(#) = qualifier out of range (m) = manual integration

Quantitation Report (Not Reviewed)

Data File : M:\CHICO\DATA\C111030\1030C17W.D Vial: 1
Acq On : 31 Oct 11 00:54 Operator: STC
Sample : Voc Std 10-30-11@1.0ug/L Inst : Chico
Misc : Water 10mLw/ IS:10-30-11 Multiplr: 1.00

Quant Time: Dec 2 11:19 2011 Quant Results File: CALLW.RES

Quant Method : M:\CHICO\DATA\C111030\CALLW.M (RTE Integrator)
Title : METHOD 8260
Last Update : Fri Dec 02 11:18:49 2011
Response via : Initial Calibration
DataAcq Meth : V8260

	Compound	R.T.	QIon	Response	Conc Unit	Qvalue
95)	Hexachlorobutadiene	25.84	223	3387	0.95661 ppb	84
96)	Naphthalene	25.94	128	24301	1.01065 ppb	97
97)	1,2,3-Trichlorobenzene	26.29	180	13947	0.94613 ppb	97

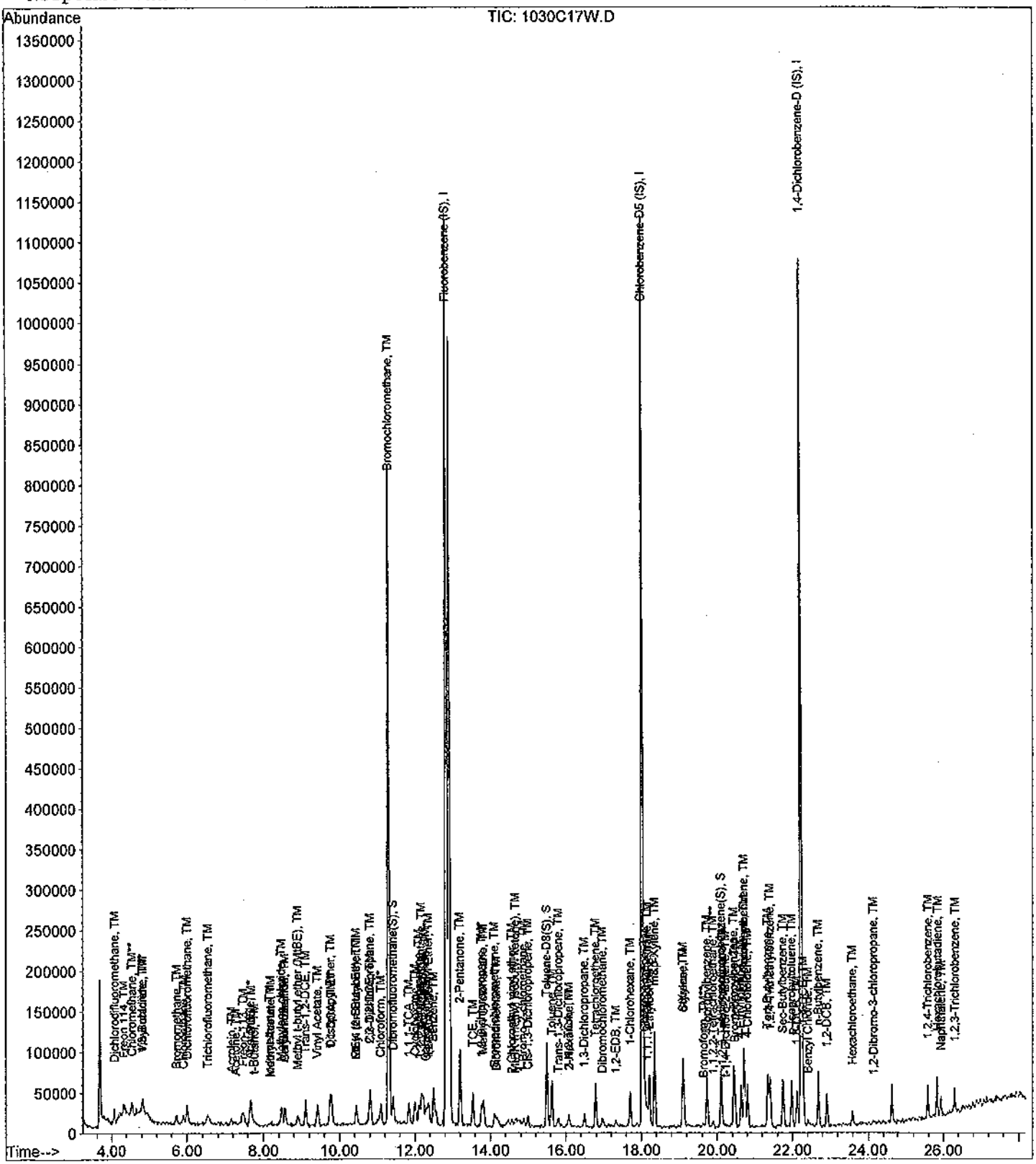
(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : M:\CHICO\DATA\C111030\1030C17W.D Vial: 1
Acq On : 31 Oct 11 00:54 Operator: STC
Sample : Voc Std 10-30-11@1.0ug/L Inst : Chico
Misc : Water 10mLw/ IS:10-30-11 Multiplr: 1.00

Quant Time: Dec 2 11:19 2011 Quant Results File: CALLW.RES

Method : M:\CHICO\DATA\C111030\CALLW.M (RTE Integrator)
Title : METHOD 8260
Last Update : Fri Dec 02 11:18:49 2011
Response via : Initial Calibration



Quantitation Report (Not Reviewed)

Data File : M:\CHICO\DATA\C111030\1030C18W.D Vial: 1
Acq On : 31 Oct 11 1:37 Operator: STC
Sample : Voc Std 10-30-11@2.0ug/L Inst : Chico
Misc : Water 10mLw/ IS:10-30-11 Multiplr: 1.00

Quant Time: Dec 2 11:19 2011 Quant Results File: CALLW.RES

Quant Method : M:\CHICO\DATA\C111030\CALLW.M (RTE Integrator)
Title : METHOD 8260
Last Update : Fri Dec 02 11:18:49 2011
Response via : Initial Calibration
DataAcq Meth : V8260

Internal Standards	R.T.	QIon	Response	Conc Units		Dev(Min)
1) Fluorobenzene (IS)	12.83	96	543693	25.00000 ppb		0.00
55) Chlorobenzene-D5 (IS)	18.04	117	392832	25.00000 ppb		0.00
71) 1,4-Dichlorobenzene-D (IS)	22.24	152	191296	25.00000 ppb		0.00
System Monitoring Compounds						
33) Dibromofluoromethane(S)	11.42	111	60078	4.14811 ppb		0.00
Spiked Amount 25.097			Recovery = 16.528%			
38) 1,2-DCA-D4(S)	12.22	65	54589	4.23414 ppb		0.00
Spiked Amount 24.225			Recovery = 17.478%			
56) Toluene-D8(S)	15.50	98	219343	3.96834 ppb		0.00
Spiked Amount 25.808			Recovery = 15.375%			
64) 4-Bromofluorobenzene(S)	20.11	95	76951	3.88546 ppb		0.00
Spiked Amount 25.459			Recovery = 15.260%			
Target Compounds						Qvalue
2) Dichlorodifluoromethane	4.07	85	36846	1.83942 ppb		95
3) Freon 114	4.32	85	25225	2.01066 ppb		95
4) Chloromethane	4.55	50	51493	2.07441 ppb		100
5) Vinyl chloride	4.81	62	31638	1.90546 ppb		92
6) 1,3-Butadiene	4.84	54	317	6.83143 ppb	#	82
7) Bromomethane	5.71	94	28199	2.33962 ppb		74
8) Chloroethane	5.92	64	27665	2.01714 ppb		90
9) Dichlorofluoromethane	6.00	67	80153	2.11371 ppb		99
10) Trichlorofluoromethane	6.54	101	46217	2.05380 ppb		88
11) Acetonitrile	7.66	41	44013	73.89637 ug/l		100
12) Acrolein	7.16	56	19054	69.90947 ppb		98
13) Acetone	7.27	43	7499	4.81405 ppb	#	42
14) Freon-113	7.46	101	27051	1.13214 ppb		88
15) 1,1-DCE	7.68	96	30774	1.98259 ppb		95
16) t-Butanol	7.77	59	6037	81.90616 ppb	#	90
17) Methyl Acetate	8.19	43	14018	2.50439 ppb		92
18) Iodomethane	8.16	142	9850	4.55459 ppb	#	90
19) Acrylonitrile	8.55	53	4257	2.21042 ppb	#	19
20) Methylene chloride	8.48	84	31366	2.11859 ppb		85
21) Carbon disulfide	8.57	76	31256	2.07243 ppb		98
22) Methyl t-butyl ether (MtBE	8.91	73	49153	2.09506 ppb		96
23) Trans-1,2-DCE	9.10	96	38306	2.41590 ppb		94
24) Diisopropyl Ether	9.76	45	107058	2.06409 ppb		95
25) 1,1-DCA	9.80	63	63745	2.07237 ppb		98
26) Vinyl Acetate	9.41	43	25227	1.48278 ppb		94
27) Ethyl tert Butyl Ether	10.45	59	76607	2.16342 ppb		92
28) MEK (2-Butanone)	10.43	43	15669	2.02476 ppb		99
29) Cis-1,2-DCE	10.80	96	37483	2.02559 ppb		88
30) 2,2-Dichloropropane	10.81	77	45572	2.06826 ppb		95
31) Chloroform	11.08	83	60802	2.05473 ppb		96
32) Bromochloromethane	11.31	128	10121	1.96459 ppb		82
34) 1,1,1-TCA	11.83	97	53989	2.00623 ppb		93
35) Cyclohexane	12.00	56	52880	2.11050 ppb		95
36) 1,1-Dichloropropene	12.11	75	46149	2.00137 ppb		96
37) 2,2,4-Trimethylpentane	12.19	57	95598	1.55463 ppb		92
39) Carbon Tetrachloride	12.30	117	36525	1.97107 ppb		97

(#) = qualifier out of range (m) = manual integration

Quantitation Report (Not Reviewed)

Data File : M:\CHICO\DATA\C111030\1030C18W.D Vial: 1
Acq On : 31 Oct 11 1:37 Operator: STC
Sample : Voc Std 10-30-11@2.0ug/L Inst : Chico
Misc : Water 10mLw/ IS:10-30-11 Multiplr: 1.00

Quant Time: Dec 2 11:19 2011 Quant Results File: CALLW.RES

Quant Method : M:\CHICO\DATA\C111030\CALLW.M (RTE Integrator)
Title : METHOD 8260
Last Update : Fri Dec 02 11:18:49 2011
Response via : Initial Calibration
DataAcq Meth : V8260

Compound	R.T.	QIon	Response	Conc Unit		Qvalue
40) Tert Amyl Methyl Ether	12.36	73	55233	2.08607 ppb	#	90
41) 1,2-DCA	12.38	62	31308	2.06118 ppb	#	89
42) Benzene	12.50	78	136536	2.06095 ppb		98
43) TCE	13.53	95	39821	2.17034 ppb		94
44) 2-Pentanone	13.20	43	299963	78.15618 ppb		98
45) 1,2-Dichloropropane	13.76	63	28619	1.90057 ppb		96
46) Bromodichloromethane	14.11	83	33573	1.95155 ppb		94
47) Methyl Cyclohexane	13.82	83	40581	1.89267 ppb		83
48) Dibromomethane	14.18	93	12379	2.05557 ppb		85
49) 2-Chloroethyl vinyl ether	14.57	63	8433	2.20290 ppb	#	78
50) 1-Bromo-2-chloroethane	14.88	63	26134	2.03401 ppb	#	77
51) Cis-1,3-Dichloropropene	15.00	75	32272	1.96729 ppb		92
52) Toluene	15.64	91	132018	2.02040 ppb		99
53) Trans-1,3-Dichloropropene	15.80	75	23617	1.99976 ppb		93
54) 1,1,2-TCA	16.07	83	13091	2.05665 ppb		90
57) 1,2-EDB	17.33	107	13384	1.79015 ppb	#	92
58) Tetrachloroethene	16.79	164	40109	1.98595 ppb		92
59) 1-Chlorohexane	17.70	91	46192	1.98613 ppb		96
60) 1,1,1,2-Tetrachloroethane	18.15	131	23520	1.86019 ppb		89
61) m&p-Xylene	18.36	106	110324	3.69766 ppb		95
62) o-Xylene	19.09	106	52936	1.84480 ppb		88
63) Styrene	19.12	104	82149	1.89672 ppb		97
65) 2-Hexanone	16.12	43	7490	2.08373 ppb		85
66) 1,3-Dichloropropane	16.50	76	29140	1.97652 ppb		91
67) Dibromochloromethane	16.96	129	16360	1.69982 ppb		100
68) Chlorobenzene	18.11	112	79768	1.86890 ppb		98
69) Ethylbenzene	18.21	91	149915	1.88621 ppb		93
70) Bromoform	19.64	173	6510	2.30500 ppb	#	44
72) MIBK (methyl isobutyl keto	14.67	43	12611	2.31897 ppb	#	76
73) Isopropylbenzene	19.73	105	142568	2.05630 ppb		95
74) 1,1,2,2-Tetrachloroethane	19.90	83	12016	2.07037 ppb	#	84
75) 1,2,3-Trichloropropane	20.16	110	1264	2.01509 ppb	#	74
76) t-1,4-Dichloro-2-Butene	20.21	53	2422	1.84044 ppb	#	55
77) Bromobenzene	20.48	156	33262	2.07982 ppb		93
78) n-Propylbenzene	20.45	91	180141	2.17796 ppb		94
79) 4-Ethyltoluene	20.64	105	114833	2.00627 ppb		95
80) 2-Chlorotoluene	20.74	91	114105	2.08302 ppb		94
81) 1,3,5-Trimethylbenzene	20.72	105	116458	2.06813 ppb		96
82) 4-Chlorotoluene	20.81	91	100754	2.13600 ppb		94
83) Tert-Butylbenzene	21.36	119	127189	2.08627 ppb		98
84) 1,2,4-Trimethylbenzene	21.41	105	117571	1.99911 ppb		93
85) Sec-Butylbenzene	21.75	105	153012	2.09286 ppb		92
86) p-Isopropyltoluene	21.99	119	133490	2.13166 ppb		95
87) Benzyl Chloride	22.42	91	15798	1.90043 ppb	#	85
88) 1,3-DCB	22.13	146	68550	2.09596 ppb		97
89) 1,4-DCB	22.29	146	62433	2.05672 ppb		94
90) Hexachloroethane	23.60	117	11613	2.93003 ppb		93
91) n-Butylbenzene	22.70	91	110023	2.01433 ppb		97
92) 1,2-DCB	22.93	146	54129	2.08070 ppb		92
93) 1,2-Dibromo-3-chloropropan	24.14	155	1335	2.54260 ppb		84
94) 1,2,4-Trichlorobenzene	25.58	180	37932	2.01219 ppb		89

(#) = qualifier out of range (m) = manual integration

Quantitation Report (Not Reviewed)

Data File : M:\CHICO\DATA\C111030\1030C18W.D Vial: 1
Acq On : 31 Oct 11 1:37 Operator: STC
Sample : Voc Std 10-30-11@2.0ug/L Inst : Chico
Misc : Water 10mLw/ IS:10-30-11 Multiplr: 1.00

Quant Time: Dec 2 11:19 2011 Quant Results File: CALLW.RES

Quant Method : M:\CHICO\DATA\C111030\CALLW.M (RTE Integrator)
Title : METHOD 8260
Last Update : Fri Dec 02 11:18:49 2011
Response via : Initial Calibration
DataAcq Meth : V8260

Compound	R.T.	QIon	Response	Conc	Unit		Qvalue
95) Hexachlorobutadiene	25.84	223	6417	1.87363	ppb	#	88
96) Naphthalene	25.93	128	46421	1.99582	ppb		98
97) 1,2,3-Trichlorobenzene	26.30	180	33171	2.32627	ppb		93

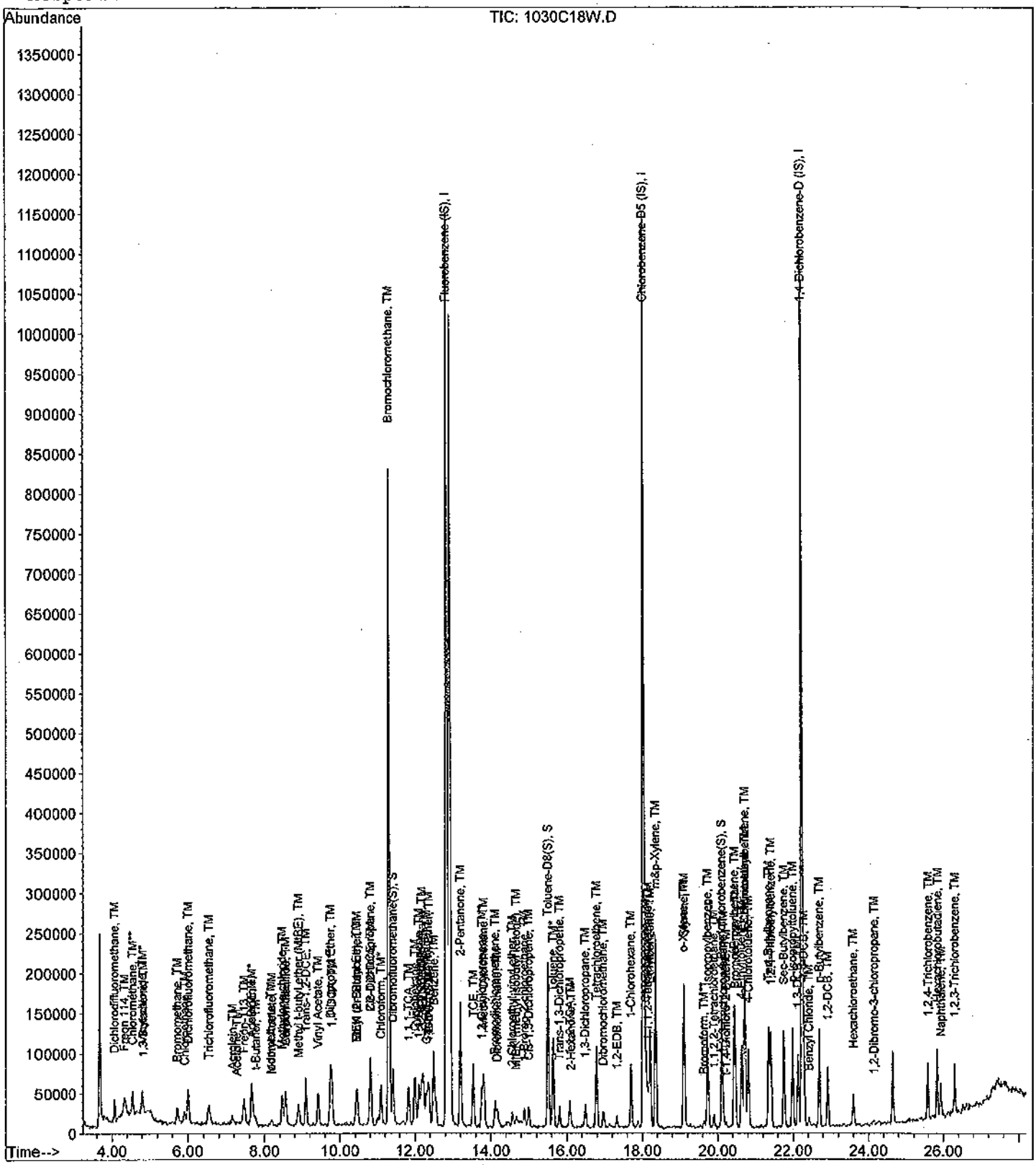
(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : M:\CHICO\DATA\C111030\1030C18W.D Vial: 1
Acq On : 31 Oct 11 1:37 Operator: STC
Sample : Voc Std 10-30-11@2.0ug/L Inst : Chico
Misc : Water 10mLw/ IS:10-30-11 Multiplr: 1.00

Quant Time: Dec 2 11:19 2011 Quant Results File: CALLW.RES

Method : M:\CHICO\DATA\C111030\CALLW.M (RTE Integrator)
Title : METHOD 8260
Last Update : Fri Dec 02 11:18:49 2011
Response via : Initial Calibration



Quantitation Report (Not Reviewed)

Data File : M:\CHICO\DATA\C111030\1030C19W.D Vial: 1
Acq On : 31 Oct 11 2:20 Operator: STC
Sample : Voc Std 10-30-11@5.0ug/L Inst : Chico
Misc : Water 10mLw/ IS:10-30-11 Multiplr: 1.00

Quant Time: Dec 2 11:19 2011 Quant Results File: CALLW.RES

Quant Method : M:\CHICO\DATA\C111030\CALLW.M (RTE Integrator)
Title : METHOD 8260
Last Update : Fri Dec 02 11:18:49 2011
Response via : Initial Calibration
DataAcq Meth : V8260

Internal Standards	R.T.	QIon	Response	Conc	Units		Dev(Min)
1) Fluorobenzene (IS)	12.84	96	541888	25.00000	ppb		0.00
55) Chlorobenzene-D5 (IS)	18.03	117	369024	25.00000	ppb		0.00
71) 1,4-Dichlorobenzene-D (IS)	22.24	152	201600	25.00000	ppb		0.00
System Monitoring Compounds							
33) Dibromofluoromethane(S)	11.43	111	148052	10.25637	ppb		0.00
Spiked Amount 25.097			Recovery = 40.865%				
38) 1,2-DCA-D4(S)	12.23	65	131632	10.24391	ppb		0.00
Spiked Amount 24.225			Recovery = 42.286%				
56) Toluene-D8(S)	15.50	98	537545	10.35266	ppb		0.00
Spiked Amount 25.808			Recovery = 40.115%				
64) 4-Bromofluorobenzene(S)	20.11	95	187725	10.09026	ppb		0.00
Spiked Amount 25.459			Recovery = 39.632%				
Target Compounds							Qvalue
2) Dichlorodifluoromethane	4.07	85	95693	4.79307	ppb		98
3) Freon 114	4.35	85	60331	4.82494	ppb		99
4) Chloromethane	4.55	50	115016	4.64888	ppb		98
5) Vinyl chloride	4.82	62	95476	5.76939	ppb		92
6) 1,3-Butadiene	4.78	54	267	5.77308	ppb	#	1
7) Bromomethane	5.73	94	55608	4.62906	ppb		89
8) Chloroethane	5.91	64	65920	4.82244	ppb		93
9) Dichlorofluoromethane	6.01	67	193619	5.12293	ppb		100
10) Trichlorofluoromethane	6.52	101	112493	5.01564	ppb		99
11) Acetonitrile	7.65	41	62470	105.23443	ug/l		100
12) Acrolein	7.15	56	26911	99.06583	ppb		94
13) Acetone	7.29	43	12573	8.09824	ppb		97
14) Freon-113	7.46	101	68643	4.67447	ppb		88
15) 1,1-DCE	7.67	96	76091	4.91843	ppb		89
16) t-Butanol	7.76	59	7217	98.24180	ppb		93
17) Methyl Acetate	8.19	43	20340	3.95990	ppb		95
18) Iodomethane	8.17	142	26092	5.76882	ppb		90
19) Acrylonitrile	8.56	53	8927	5.07063	ppb		92
20) Methylene chloride	8.48	84	71913	4.87349	ppb		94
21) Carbon disulfide	8.56	76	72344	4.81275	ppb		99
22) Methyl t-butyl ether (MtBE	8.90	73	114808	4.90979	ppb	#	90
23) Trans-1,2-DCE	9.10	96	89303	5.65097	ppb		95
24) Diisopropyl Ether	9.75	45	260084	5.03115	ppb		91
25) 1,1-DCA	9.79	63	163680	5.33900	ppb		96
26) Vinyl Acetate	9.42	43	54231	4.98892	ppb	#	79
27) Ethyl tert Butyl Ether	10.44	59	180946	5.12702	ppb		98
28) MEK (2-Butanone)	10.43	43	31144	4.71582	ppb	#	87
29) Cis-1,2-DCE	10.82	96	90101	4.88531	ppb		94
30) 2,2-Dichloropropane	10.82	77	110498	5.03159	ppb		99
31) Chloroform	11.10	83	153864	5.21698	ppb		100
32) Bromochloromethane	11.31	128	28344	5.52019	ppb		86
34) 1,1,1-TCA	11.84	97	133883	4.99165	ppb		96
35) Cyclohexane	12.00	56	119918	4.80199	ppb		98
36) 1,1-Dichloropropene	12.10	75	117037	5.09252	ppb		98
37) 2,2,4-Trimethylpentane	12.18	57	209064	4.78823	ppb		95
39) Carbon Tetrachloride	12.30	117	94473	5.11522	ppb		93

(#) = qualifier out of range (m) = manual integration

Quantitation Report (Not Reviewed)

Data File : M:\CHICO\DATA\C111030\1030C19W.D Vial: 1
Acq On : 31 Oct 11 2:20 Operator: STC
Sample : Voc Std 10-30-11@5.0ug/L Inst : Chico
Misc : Water 10mLw/ IS:10-30-11 Multiplr: 1.00

Quant Time: Dec 2 11:19 2011 Quant Results File: CALLW.RES

Quant Method : M:\CHICO\DATA\C111030\CALLW.M (RTE Integrator)
Title : METHOD 8260
Last Update : Fri Dec 02 11:18:49 2011
Response via : Initial Calibration
DataAcq Meth : V8260

Compound	R.T.	QIon	Response	Conc Unit		Qvalue
40) Tert Amyl Methyl Ether	12.36	73	128972	4.88731 ppb		98
41) 1,2-DCA	12.38	62	76954	5.08320 ppb		98
42) Benzene	12.50	78	331433	5.01951 ppb		100
43) TCE	13.54	95	94771	5.18246 ppb		92
44) 2-Pentanone	13.20	43	389562	101.83954 ppb		98
45) 1,2-Dichloropropane	13.77	63	79197	5.27695 ppb		96
46) Bromodichloromethane	14.12	83	87445	5.09999 ppb		96
47) Methyl Cyclohexane	13.82	83	103751	4.85498 ppb		94
48) Dibromomethane	14.16	93	31588	5.26275 ppb		82
49) 2-Chloroethyl vinyl ether	14.58	63	20703	5.42613 ppb	#	92
50) 1-Bromo-2-chloroethane	14.88	63	66427	5.18724 ppb		82
51) Cis-1,3-Dichloropropene	15.01	75	82661	5.05577 ppb		98
52) Toluene	15.63	91	327752	5.03263 ppb		100
53) Trans-1,3-Dichloropropene	15.80	75	58149	4.94014 ppb		92
54) 1,1,2-TCA	16.08	83	31919	5.03132 ppb		92
57) 1,2-EDB	17.32	107	37026	5.27183 ppb	#	70
58) Tetrachloroethene	16.79	164	96522	5.08752 ppb		93
59) 1-Chlorohexane	17.70	91	112708	5.15879 ppb		93
60) 1,1,1,2-Tetrachloroethane	18.16	131	63531	5.34881 ppb		97
61) m&p-Xylene	18.36	106	280238	9.99854 ppb		95
62) o-Xylene	19.11	106	143606	5.32749 ppb		99
63) Styrene	19.12	104	215758	5.30297 ppb		96
65) 2-Hexanone	16.11	43	15947	4.72270 ppb		87
66) 1,3-Dichloropropane	16.50	76	70642	5.10067 ppb		92
67) Dibromochloromethane	16.97	129	44643	4.93770 ppb		89
68) Chlorobenzene	18.10	112	203928	5.08611 ppb		100
69) Ethylbenzene	18.22	91	385898	5.16857 ppb		100
70) Bromoform	19.63	173	18516	4.67993 ppb		89
72) MIBK (methyl isobutyl keto	14.68	43	24562	4.28572 ppb		100
73) Isopropylbenzene	19.74	105	367532	5.03008 ppb		99
74) 1,1,2,2-Tetrachloroethane	19.90	83	28865	4.71928 ppb	#	80
75) 1,2,3-Trichloropropane	20.15	110	3527	5.18588 ppb	#	55
76) t-1,4-Dichloro-2-Butene	20.22	53	6664	4.80505 ppb	#	82
77) Bromobenzene	20.48	156	83579	4.95894 ppb		96
78) n-Propylbenzene	20.44	91	444891	5.10396 ppb		99
79) 4-Ethyltoluene	20.64	105	295868	4.90496 ppb		95
80) 2-Chlorotoluene	20.73	91	295014	5.11030 ppb		95
81) 1,3,5-Trimethylbenzene	20.71	105	303247	5.11000 ppb		98
82) 4-Chlorotoluene	20.82	91	244634	4.92119 ppb		94
83) Tert-Butylbenzene	21.36	119	322939	5.02640 ppb		96
84) 1,2,4-Trimethylbenzene	21.42	105	306527	4.94563 ppb		98
85) Sec-Butylbenzene	21.76	105	387689	5.03168 ppb		98
86) p-Isopropyltoluene	21.99	119	333281	5.05005 ppb		97
87) Benzyl Chloride	22.43	91	41624	4.75127 ppb		92
88) 1,3-DCB	22.14	146	173843	5.04369 ppb		91
89) 1,4-DCB	22.30	146	155968	4.87542 ppb		94
90) Hexachloroethane	23.60	117	36490	4.84597 ppb		81
91) n-Butylbenzene	22.70	91	281770	4.89505 ppb		94
92) 1,2-DCB	22.93	146	134998	4.92405 ppb		97
93) 1,2-Dibromo-3-chloropropan	24.14	155	4220	5.04887 ppb		94
94) 1,2,4-Trichlorobenzene	25.59	180	99085	4.98753 ppb		96

(#) = qualifier out of range (m) = manual integration

Quantitation Report (Not Reviewed)

Data File : M:\CHICO\DATA\C111030\1030C19W.D Vial: 1
Acq On : 31 Oct 11 2:20 Operator: STC
Sample : Voc Std 10-30-11@5.0ug/L Inst : Chico
Misc : Water 10mLw/ IS:10-30-11 Multiplr: 1.00

Quant Time: Dec 2 11:19 2011 Quant Results File: CALLW.RES

Quant Method : M:\CHICO\DATA\C111030\CALLW.M (RTE Integrator)
Title : METHOD 8260
Last Update : Fri Dec 02 11:18:49 2011
Response via : Initial Calibration
DataAcq Meth : V8260

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
95) Hexachlorobutadiene	25.84	223	15832	4.38634	ppb	86
96) Naphthalene	25.94	128	114277	4.66210	ppb	98
97) 1,2,3-Trichlorobenzene	26.29	180	73429	4.88636	ppb	92

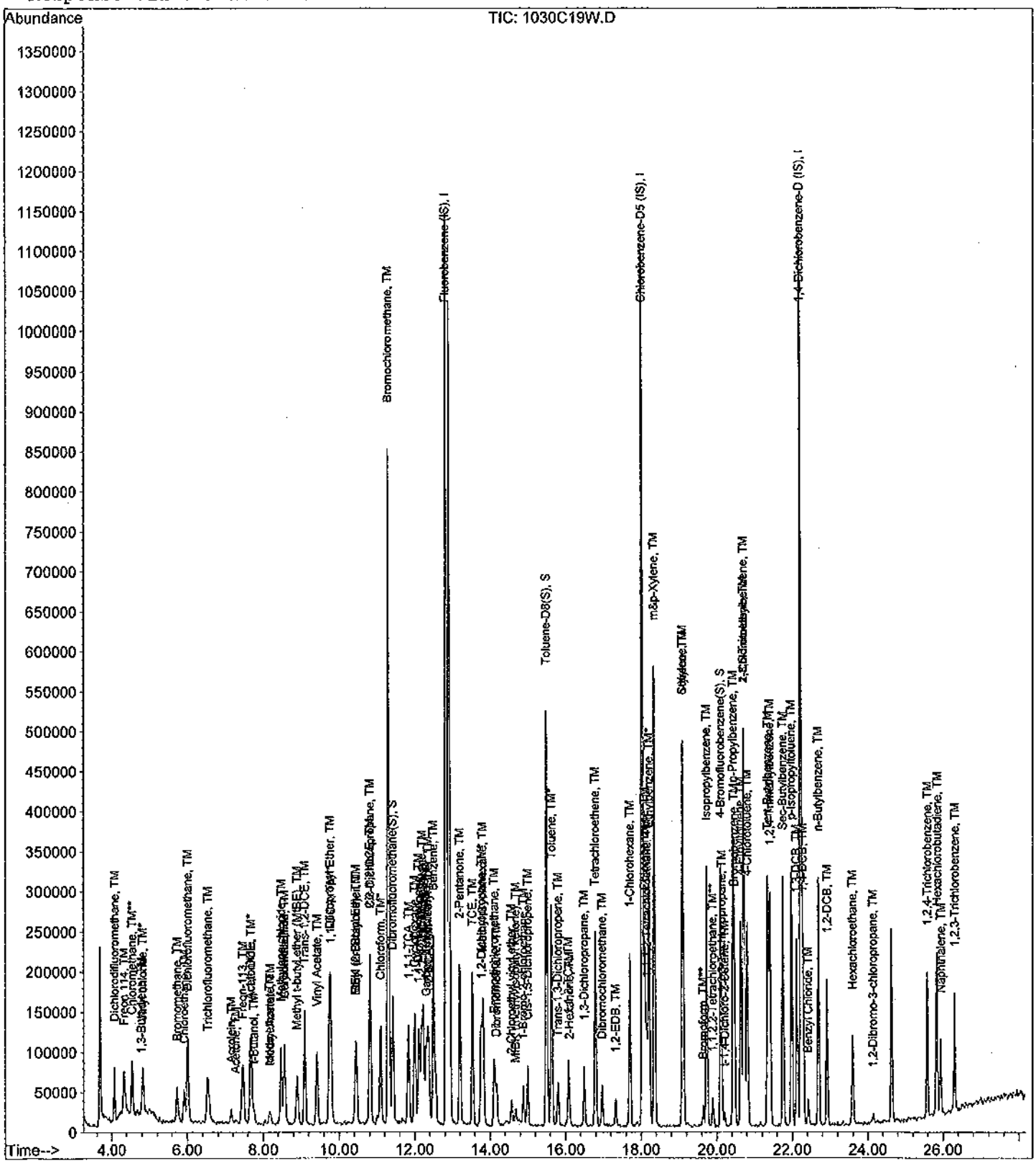
(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : M:\CHICO\DATA\C111030\1030C19W.D Vial: 1
Acq On : 31 Oct 11 2:20 Operator: STC
Sample : Voc Std 10-30-11@5.0ug/L Inst : Chico
Misc : Water 10mLw/ IS:10-30-11 Multiplr: 1.00

Quant Time: Dec 2 11:19 2011 Quant Results File: CALLW.RES

Method : M:\CHICO\DATA\C111030\CALLW.M (RTE Integrator)
Title : METHOD 8260
Last Update : Fri Dec 02 11:18:49 2011
Response via : Initial Calibration



Quantitation Report (Not Reviewed)

Data File : M:\CHICO\DATA\C111030\1030C20W.D Vial: 1
Acq On : 31 Oct 11 3:03 Operator: STC
Sample : Voc Std 10-30-11@10ug/L Inst : Chico
Misc : Water 10mLw/ IS:10-30-11 Multiplr: 1.00

Quant Time: Dec 2 11:19 2011 Quant Results File: CALLW.RES

Quant Method : M:\CHICO\DATA\C111030\CALLW.M (RTE Integrator)
Title : METHOD 8260
Last Update : Fri Dec 02 11:18:49 2011
Response via : Initial Calibration
DataAcq Meth : V8260

Internal Standards	R.T.	QIon	Response	Conc Units	Dev(Min)
1) Fluorobenzene (IS)	12.84	96	556544	25.00000 ppb	0.00
55) Chlorobenzene-D5 (IS)	18.04	117	375296	25.00000 ppb	0.00
71) 1,4-Dichlorobenzene-D (IS)	22.24	152	203520	25.00000 ppb	0.00
System Monitoring Compounds					
33) Dibromofluoromethane(S)	11.43	111	379665	25.60885 ppb	0.00
Spiked Amount 25.097			Recovery = 102.039%		
38) 1,2-DCA-D4(S)	12.23	65	325575	24.66979 ppb	0.00
Spiked Amount 24.225			Recovery = 101.836%		
56) Toluene-D8(S)	15.50	98	1312175	24.84903 ppb	0.00
Spiked Amount 25.808			Recovery = 96.283%		
64) 4-Bromofluorobenzene(S)	20.11	95	480879	25.41539 ppb	0.00
Spiked Amount 25.459			Recovery = 99.825%		
Target Compounds					Qvalue
2) Dichlorodifluoromethane	4.07	85	203737	9.93605 ppb	100
3) Freon 114	4.33	85	139352	10.85112 ppb	100
4) Chloromethane	4.55	50	239697	9.43328 ppb	100
5) Vinyl chloride	4.82	62	177123	10.42127 ppb	100
6) 1,3-Butadiene	4.83	54	475	10.00000 ppb	100
7) Bromomethane	5.73	94	118673	9.61873 ppb	100
8) Chloroethane	5.91	64	132559	9.44211 ppb	100
9) Dichlorofluoromethane	6.01	67	384006	9.89278 ppb	100
10) Trichlorofluoromethane	6.53	101	230965	10.02667 ppb	100
11) Acetonitrile	7.65	41	74340	121.93234 ug/l	100
12) Acrolein	7.16	56	32973	118.18505 ppb	100
13) Acetone	7.28	43	21604	13.54865 ppb	100
14) Freon-113	7.47	101	140200	10.44199 ppb	100
15) 1,1-DCE	7.68	96	152367	9.58945 ppb	100
16) t-Butanol	7.77	59	9738	129.06824 ppb	100
17) Methyl Acetate	8.19	43	49512	10.32908 ppb	100
18) Iodomethane	8.16	142	66421	8.64718 ppb	100
19) Acrylonitrile	8.56	53	18948	10.88495 ppb	100
20) Methylene chloride	8.47	84	148076	9.77073 ppb	100
21) Carbon disulfide	8.56	76	151616	9.82077 ppb	100
22) Methyl t-butyl ether (MtBE	8.89	73	244998	10.20148 ppb	100
23) Trans-1,2-DCE	9.10	96	182548	11.24718 ppb	100
24) Diisopropyl Ether	9.76	45	539900	10.16896 ppb	100
25) 1,1-DCA	9.79	63	326209	10.36025 ppb	100
26) Vinyl Acetate	9.42	43	98410	10.00173 ppb	100
27) Ethyl tert Butyl Ether	10.45	59	389708	10.75141 ppb	100
28) MEK (2-Butanone)	10.44	43	65986	10.45330 ppb	100
29) Cis-1,2-DCE	10.82	96	187720	9.91021 ppb	100
30) 2,2-Dichloropropane	10.81	77	219771	9.74386 ppb	100
31) Chloroform	11.09	83	311468	10.28266 ppb	100
32) Bromochloromethane	11.32	128	54568	10.34763 ppb	100
34) 1,1,1-TCA	11.84	97	285282	10.35627 ppb	100
35) Cyclohexane	12.00	56	254779	9.93369 ppb	100
36) 1,1-Dichloropropene	12.11	75	236871	10.03533 ppb	100
37) 2,2,4-Trimethylpentane	12.18	57	404324	10.03470 ppb	100
39) Carbon Tetrachloride	12.31	117	201043	10.59877 ppb	100

(#) = qualifier out of range (m) = manual integration

Quantitation Report (Not Reviewed)

Data File : M:\CHICO\DATA\C111030\1030C20W.D Vial: 1
Acq On : 31 Oct 11 3:03 Operator: STC
Sample : Voc Std 10-30-11@10ug/L Inst : Chico
Misc : Water 10mLw/ IS:10-30-11 Multiplr: 1.00

Quant Time: Dec 2 11:19 2011 Quant Results File: CALLW.RES

Quant Method : M:\CHICO\DATA\C111030\CALLW.M (RTE Integrator)
Title : METHOD 8260
Last Update : Fri Dec 02 11:18:49 2011
Response via : Initial Calibration
DataAcq Meth : V8260

Compound	R.T.	QIon	Response	Conc Unit	Qvalue
40) Tert Amyl Methyl Ether	12.34	73	277234	10.22896 ppb	100
41) 1,2-DCA	12.38	62	162875	10.47539 ppb	100
42) Benzene	12.50	78	685232	10.10446 ppb	100
43) TCE	13.53	95	193489	10.30213 ppb	100
44) 2-Pentanone	13.20	43	514551	130.97195 ppb	100
45) 1,2-Dichloropropane	13.77	63	160165	10.39087 ppb	100
46) Bromodichloromethane	14.11	83	192788	10.94774 ppb	100
47) Methyl Cyclohexane	13.81	83	223185	10.16882 ppb	100
48) Dibromomethane	14.17	93	67961	11.02453 ppb	100
49) 2-Chloroethyl vinyl ether	14.58	63	41287	10.53611 ppb	100
50) 1-Bromo-2-chloroethane	14.89	63	134227	10.20568 ppb	100
51) Cis-1,3-Dichloropropene	15.00	75	178053	10.60342 ppb	100
52) Toluene	15.63	91	682617	10.20555 ppb	100
53) Trans-1,3-Dichloropropene	15.80	75	130192	10.76940 ppb	100
54) 1,1,2-TCA	16.08	83	73187	11.23251 ppb	100
57) 1,2-EDB	17.33	107	76477	10.70696 ppb	100
58) Tetrachloroethene	16.79	164	188693	9.77949 ppb	100
59) 1-Chlorohexane	17.70	91	223919	10.07778 ppb	100
60) 1,1,1,2-Tetrachloroethane	18.16	131	136098	11.26688 ppb	100
61) m&p-Xylene	18.36	106	573288	20.11236 ppb	100
62) o-Xylene	19.11	106	288268	10.51543 ppb	100
63) Styrene	19.13	104	435830	10.53295 ppb	100
65) 2-Hexanone	16.11	43	35479	10.33151 ppb	100
66) 1,3-Dichloropropane	16.49	76	155699	11.05429 ppb	100
67) Dibromochloromethane	16.98	129	100610	10.94190 ppb	100
68) Chlorobenzene	18.11	112	417306	10.23397 ppb	100
69) Ethylbenzene	18.22	91	783451	10.31788 ppb	100
70) Bromoform	19.64	173	44444	9.50336 ppb	100
72) MIBK (methyl isobutyl keto	14.68	43	56876	9.83043 ppb	100
73) Isopropylbenzene	19.73	105	751023	10.18160 ppb	100
74) 1,1,2,2-Tetrachloroethane	19.90	83	68052	11.02118 ppb	100
75) 1,2,3-Trichloropropane	20.16	110	8339	12.02369 ppb	100
76) t-1,4-Dichloro-2-Butene	20.24	53	14863	10.61580 ppb	100
77) Bromobenzene	20.48	156	169233	9.94627 ppb	100
78) n-Propylbenzene	20.44	91	904419	10.27795 ppb	100
79) 4-Ethyltoluene	20.63	105	616295	10.12067 ppb	100
80) 2-Chlorotoluene	20.74	91	594233	10.19632 ppb	100
81) 1,3,5-Trimethylbenzene	20.72	105	587753	9.81075 ppb	100
82) 4-Chlorotoluene	20.82	91	501553	9.99433 ppb	100
83) Tert-Butylbenzene	21.36	119	667298	10.28820 ppb	100
84) 1,2,4-Trimethylbenzene	21.42	105	611300	9.76991 ppb	100
85) Sec-Butylbenzene	21.76	105	815062	10.47861 ppb	100
86) p-Isopropyltoluene	21.99	119	683802	10.26358 ppb	100
87) Benzyl Chloride	22.42	91	84140	9.51375 ppb	100
88) 1,3-DCB	22.12	146	342186	9.83414 ppb	100
89) 1,4-DCB	22.30	146	328879	10.18349 ppb	100
90) Hexachloroethane	23.60	117	91222	9.10196 ppb	100
91) n-Butylbenzene	22.69	91	572922	9.85919 ppb	100
92) 1,2-DCB	22.93	146	290055	10.47995 ppb	100
93) 1,2-Dibromo-3-chloropropan	24.15	155	11552	11.48408 ppb	100
94) 1,2,4-Trichlorobenzene	25.59	180	201946	10.06924 ppb	100

(#) = qualifier out of range (m) = manual integration

Quantitation Report (Not Reviewed)

Data File : M:\CHICO\DATA\C111030\1030C20W.D Vial: 1
Acq On : 31 Oct 11 3:03 Operator: STC
Sample : Voc Std 10-30-11@10ug/L Inst : Chico
Misc : Water 10mLw/ IS:10-30-11 Multiplr: 1.00

Quant Time: Dec 2 11:19 2011 Quant Results File: CALLW.RES

Quant Method : M:\CHICO\DATA\C111030\CALLW.M (RTE Integrator)
Title : METHOD 8260
Last Update : Fri Dec 02 11:18:49 2011
Response via : Initial Calibration
DataAcq Meth : V8260

	Compound	R.T.	QIon	Response	Conc Unit		Qvalue
95)	Hexachlorobutadiene	25.84	223	37504	10.29265	ppb	100
96)	Naphthalene	25.94	128	255426	10.32217	ppb	100
97)	1,2,3-Trichlorobenzene	26.29	180	158877	10.47278	ppb	100

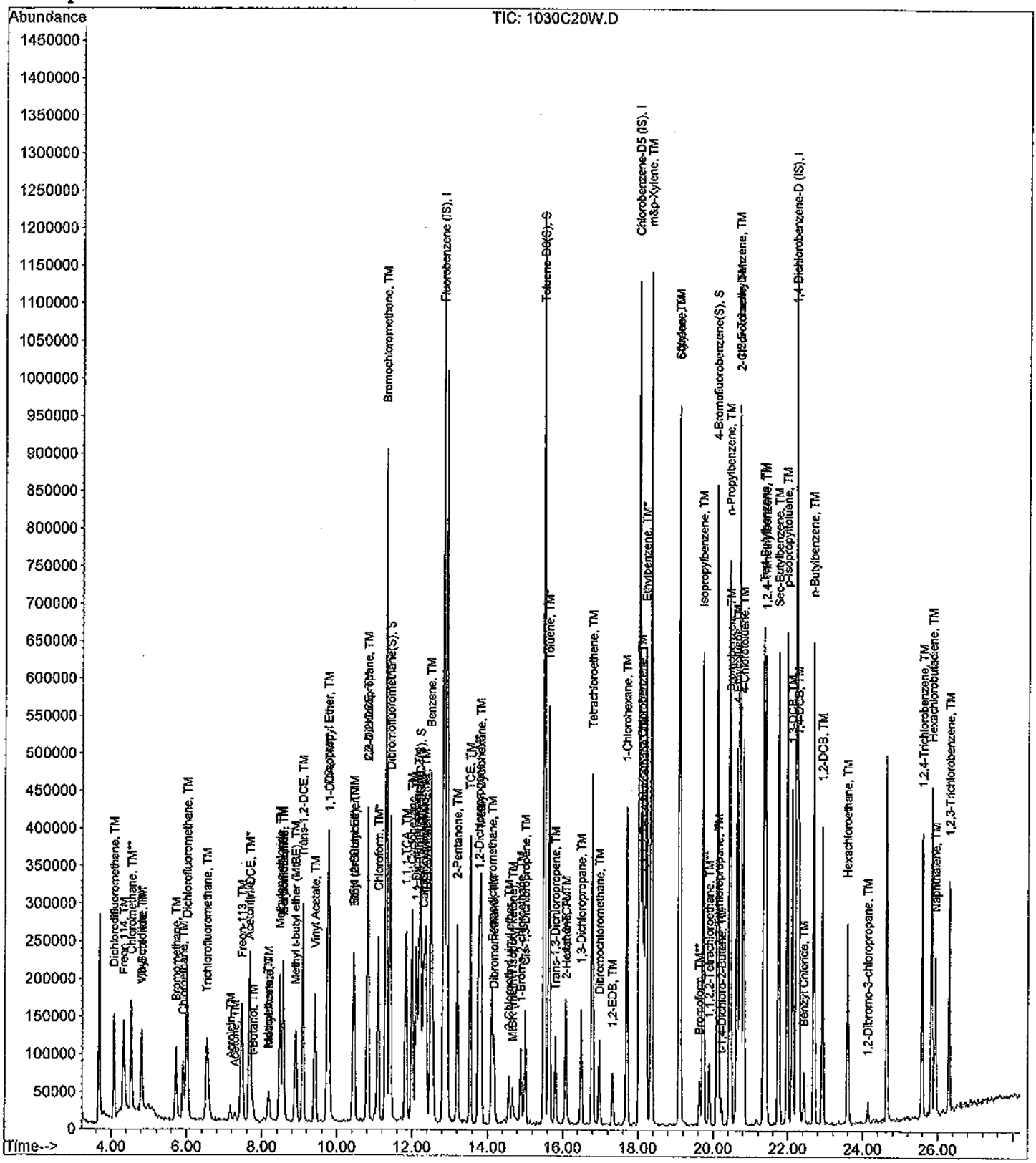
(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : M:\CHICO\DATA\C111030\1030C20W.D Vial: 1
Acq On : 31 Oct 11 3:03 Operator: STC
Sample : Voc Std 10-30-11@10ug/L Inst : Chico
Misc : Water 10mLw/ IS:10-30-11 Multiplr: 1.00

Quant Time: Dec 2 11:19 2011 Quant Results File: CALLW.RES

Method : M:\CHICO\DATA\C111030\CALLW.M (RTE Integrator)
Title : METHOD 8260
Last Update : Fri Dec 02 11:18:49 2011
Response via : Initial Calibration



Quantitation Report (Not Reviewed)

Data File : M:\CHICO\DATA\C111030\1030C21W.D Vial: 1
Acq On : 31 Oct 11 3:46 Operator: STC
Sample : Voc Std 10-30-11@20ug/L Inst : Chico
Misc : Water 10mLw/ IS:10-30-11 Multiplr: 1.00

Quant Time: Dec 2 11:19 2011 Quant Results File: CALLW.RES

Quant Method : M:\CHICO\DATA\C111030\CALLW.M (RTE Integrator)
Title : METHOD 8260
Last Update : Fri Dec 02 11:18:49 2011
Response via : Initial Calibration
DataAcq Meth : V8260

Internal Standards	R.T.	QIon	Response	Conc Units		Dev(Min)
1) Fluorobenzene (IS)	12.84	96	566784	25.00000 ppb		0.00
55) Chlorobenzene-D5 (IS)	18.03	117	371200	25.00000 ppb		0.00
71) 1,4-Dichlorobenzene-D (IS)	22.24	152	208640	25.00000 ppb		0.00
System Monitoring Compounds						
33) Dibromofluoromethane(S)	11.42	111	595137	39.41743 ppb		0.00
Spiked Amount 25.097			Recovery = 157.057%			
38) 1,2-DCA-D4(S)	12.23	65	503677	37.47559 ppb		0.00
Spiked Amount 24.225			Recovery = 154.698%			
56) Toluene-D8(S)	15.50	98	2079192	39.80872 ppb		0.00
Spiked Amount 25.808			Recovery = 154.249%			
64) 4-Bromofluorobenzene(S)	20.11	95	744294	39.77146 ppb		0.00
Spiked Amount 25.459			Recovery = 156.213%			
Target Compounds						Qvalue
2) Dichlorodifluoromethane	4.07	85	430955	20.63753 ppb		98
3) Freon 114	4.34	85	282152	21.57379 ppb		99
4) Chloromethane	4.55	50	489131	18.90198 ppb		99
5) Vinyl chloride	4.81	62	352842	20.38487 ppb		100
6) 1,3-Butadiene	4.83	54	344	7.11126 ppb	#	70
7) Bromomethane	5.73	94	258043	20.53713 ppb		86
8) Chloroethane	5.91	64	259811	18.17186 ppb		97
9) Dichlorofluoromethane	6.01	67	754464	19.08539 ppb		99
10) Trichlorofluoromethane	6.53	101	465717	19.85247 ppb		94
11) Acetonitrile	7.65	41	87890	141.55257 ug/l		100
12) Acrolein	7.16	56	38144	134.24936 ppb		98
13) Acetone	7.28	43	36583	22.52802 ppb	#	68
14) Freon-113	7.46	101	275908	21.25897 ppb		94
15) 1,1-DCE	7.68	96	292224	18.05929 ppb		95
16) t-Butanol	7.77	59	10632	138.37146 ppb		91
17) Methyl Acetate	8.19	43	90758	19.14260 ppb		98
18) Iodomethane	8.16	142	180521	16.69886 ppb	#	89
19) Acrylonitrile	8.56	53	34895	19.99122 ppb		97
20) Methylene chloride	8.48	84	296145	19.18795 ppb		91
21) Carbon disulfide	8.56	76	295040	18.76564 ppb		100
22) Methyl t-butyl ether (MtBE	8.90	73	470382	19.23240 ppb		96
23) Trans-1,2-DCE	9.11	96	343023	20.75257 ppb		92
24) Diisopropyl Ether	9.75	45	1025315	18.96280 ppb		99
25) 1,1-DCA	9.79	63	635624	19.82242 ppb		97
26) Vinyl Acetate	9.42	43	188410	20.16485 ppb		92
27) Ethyl tert Butyl Ether	10.44	59	725972	19.66653 ppb		95
28) MEK (2-Butanone)	10.44	43	124340	19.92151 ppb		98
29) Cis-1,2-DCE	10.82	96	357838	18.54986 ppb		94
30) 2,2-Dichloropropane	10.81	77	438187	19.07664 ppb		93
31) Chloroform	11.09	83	601855	19.51038 ppb		94
32) Bromochloromethane	11.32	128	108166	20.14075 ppb		85
34) 1,1,1-TCA	11.83	97	572722	20.41525 ppb		94
35) Cyclohexane	12.00	56	494926	18.94826 ppb		96
36) 1,1-Dichloropropene	12.10	75	471318	19.60721 ppb		97
37) 2,2,4-Trimethylpentane	12.18	57	798347	20.53839 ppb		93
39) Carbon Tetrachloride	12.30	117	426253	22.06560 ppb		97

(#) = qualifier out of range (m) = manual integration

Quantitation Report (Not Reviewed)

Data File : M:\CHICO\DATA\C111030\1030C21W.D Vial: 1
Acq On : 31 Oct 11 3:46 Operator: STC
Sample : Voc Std 10-30-11@20ug/L Inst : Chico
Misc : Water 10mLw/ IS:10-30-11 Multiplr: 1.00

Quant Time: Dec 2 11:19 2011 Quant Results File: CALLW.RES

Quant Method : M:\CHICO\DATA\C111030\CALLW.M (RTE Integrator)
Title : METHOD 8260
Last Update : Fri Dec 02 11:18:49 2011
Response via : Initial Calibration
DataAcq Meth : V8260

	Compound	R.T.	QIon	Response	Conc	Unit		Qvalue
40)	Tert Amyl Methyl Ether	12.35	73	534720	19.37283	ppb		97
41)	1,2-DCA	12.37	62	312083	19.70915	ppb		98
42)	Benzene	12.50	78	1326677	19.20979	ppb		99
43)	TCE	13.54	95	376706	19.69496	ppb		91
44)	2-Pentanone	13.20	43	581100	145.23880	ppb		100
45)	1,2-Dichloropropane	13.76	63	307944	19.61725	ppb		95
46)	Bromodichloromethane	14.12	83	371635	20.72255	ppb	#	93
47)	Methyl Cyclohexane	13.82	83	433011	19.37253	ppb		97
48)	Dibromomethane	14.16	93	130449	20.77893	ppb		97
49)	2-Chloroethyl vinyl ether	14.58	63	79377	19.89041	ppb		95
50)	1-Bromo-2-chloroethane	14.88	63	259762	19.39365	ppb		89
51)	Cis-1,3-Dichloropropene	15.01	75	350200	20.47835	ppb		91
52)	Toluene	15.63	91	1320740	19.38915	ppb		97
53)	Trans-1,3-Dichloropropene	15.80	75	249890	20.29730	ppb		96
54)	1,1,2-TCA	16.08	83	131106	19.75819	ppb		93
57)	1,2-EDB	17.33	107	153122	21.67400	ppb		94
58)	Tetrachloroethene	16.79	164	367302	19.24640	ppb		91
59)	1-Chlorohexane	17.70	91	453290	20.62604	ppb		98
60)	1,1,1,2-Tetrachloroethane	18.16	131	280249	23.45643	ppb		92
61)	m&p-Xylene	18.35	106	1173944	41.63932	ppb		100
62)	o-Xylene	19.11	106	585791	21.60426	ppb		95
63)	Styrene	19.12	104	892450	21.80633	ppb		99
65)	2-Hexanone	16.10	43	69030	20.32340	ppb		93
66)	1,3-Dichloropropane	16.50	76	287745	20.65470	ppb		99
67)	Dibromochloromethane	16.97	129	207497	22.81547	ppb		93
68)	Chlorobenzene	18.10	112	813528	20.17102	ppb		97
69)	Ethylbenzene	18.22	91	1522721	20.27519	ppb		95
70)	Bromoform	19.64	173	97001	19.60233	ppb		84
72)	MIBK (methyl isobutyl keto	14.68	43	106674	17.98505	ppb		92
73)	Isopropylbenzene	19.74	105	1516275	20.05166	ppb		98
74)	1,1,2,2-Tetrachloroethane	19.89	83	136632	21.58486	ppb		90
75)	1,2,3-Trichloropropane	20.15	110	13641	19.13172	ppb	#	42
76)	t-1,4-Dichloro-2-Butene	20.22	53	30320	21.12442	ppb	#	79
77)	Bromobenzene	20.48	156	337635	19.35675	ppb		93
78)	n-Propylbenzene	20.44	91	1817556	20.14810	ppb		97
79)	4-Ethyltoluene	20.64	105	1209221	19.37027	ppb		92
80)	2-Chlorotoluene	20.73	91	1199768	20.08139	ppb		99
81)	1,3,5-Trimethylbenzene	20.71	105	1237433	20.14832	ppb		99
82)	4-Chlorotoluene	20.82	91	1006043	19.55523	ppb		96
83)	Tert-Butylbenzene	21.36	119	1363292	20.50304	ppb		98
84)	1,2,4-Trimethylbenzene	21.42	105	1240098	19.33311	ppb		98
85)	Sec-Butylbenzene	21.76	105	1672276	20.97154	ppb		99
86)	p-Isopropyltoluene	21.99	119	1410527	20.65186	ppb		98
87)	Benzyl Chloride	22.43	91	179615	19.81078	ppb		96
88)	1,3-DCB	22.13	146	714268	20.02373	ppb		99
89)	1,4-DCB	22.30	146	661023	19.96579	ppb		96
90)	Hexachloroethane	23.60	117	220554	18.79960	ppb		91
91)	n-Butylbenzene	22.70	91	1175074	19.72516	ppb		100
92)	1,2-DCB	22.93	146	582656	20.53528	ppb		96
93)	1,2-Dibromo-3-chloropropan	24.14	155	19304	17.90735	ppb	#	68
94)	1,2,4-Trichlorobenzene	25.59	180	403670	19.63348	ppb		99

(#) = qualifier out of range (m) = manual integration

Quantitation Report (Not Reviewed)

Data File : M:\CHICO\DATA\C111030\1030C21W.D Vial: 1
Acq On : 31 Oct 11 3:46 Operator: STC
Sample : Voc Std 10-30-11@20ug/L Inst : Chico
Misc : Water 10mLw/ IS:10-30-11 Multiplr: 1.00

Quant Time: Dec 2 11:19 2011 Quant Results File: CALLW.RES

Quant Method : M:\CHICO\DATA\C111030\CALLW.M (RTE Integrator)
Title : METHOD 8260
Last Update : Fri Dec 02 11:18:49 2011
Response via : Initial Calibration
DataAcq Meth : V8260

Compound	R.T.	QIon	Response	Conc Unit	Qvalue
95) Hexachlorobutadiene	25.84	223	78688	21.06531 ppb	92
96) Naphthalene	25.94	128	505600	19.93070 ppb	99
97) 1,2,3-Trichlorobenzene	26.30	180	309459	19.89820 ppb	98

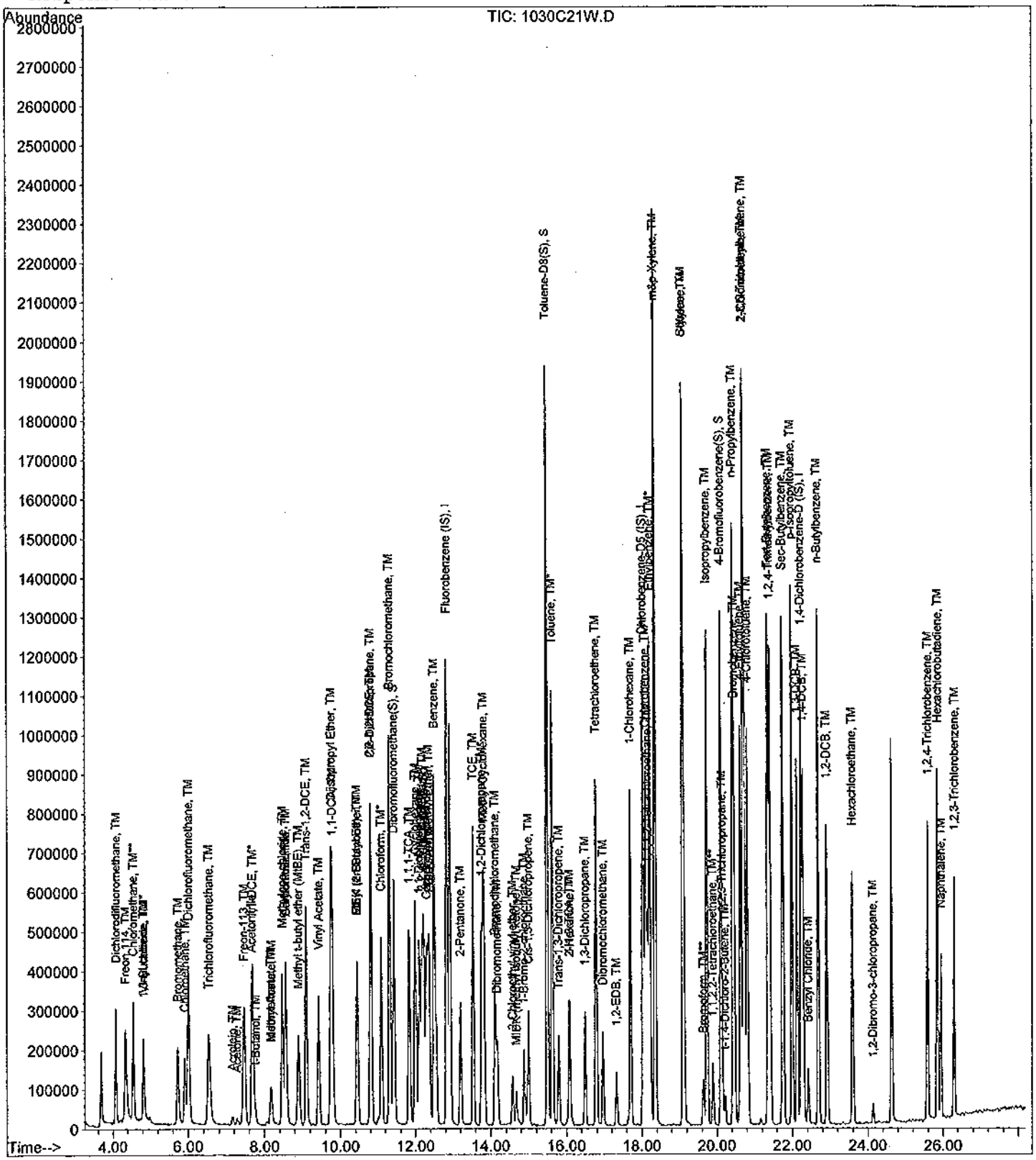
(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : M:\CHICO\DATA\C111030\1030C21W.D Vial: 1
Acq On : 31 Oct 11 3:46 Operator: STC
Sample : Voc Std 10-30-11@20ug/L Inst : Chico
Misc : Water 10mLw/ IS:10-30-11 Multiplr: 1.00

Quant Time: Dec 2 11:19 2011 Quant Results File: CALLW.RES

Method : M:\CHICO\DATA\C111030\CALLW.M (RTE Integrator)
Title : METHOD 8260
Last Update : Fri Dec 02 11:18:49 2011
Response via : Initial Calibration



Quantitation Report (Not Reviewed)

Data File : M:\CHICO\DATA\C111030\1030C22W.D Vial: 1
Acq On : 31 Oct 11 4:29 Operator: STC
Sample : Voc Std 10-30-11@40ug/L Inst : Chico
Misc : Water 10mLw/ IS:10-30-11 Multiplr: 1.00

Quant Time: Dec 2 11:19 2011 Quant Results File: CALLW.RES

Quant Method : M:\CHICO\DATA\C111030\CALLW.M (RTE Integrator)
Title : METHOD 8260
Last Update : Fri Dec 02 11:18:49 2011
Response via : Initial Calibration
DataAcq Meth : V8260

Internal Standards	R.T.	QIon	Response	Conc	Units		Dev(Min)
1) Fluorobenzene (IS)	12.84	96	576384	25.00000	ppb		0.00
55) Chlorobenzene-D5 (IS)	18.03	117	400384	25.00000	ppb		0.00
71) 1,4-Dichlorobenzene-D (IS)	22.24	152	224000	25.00000	ppb		0.00
System Monitoring Compounds							
33) Dibromofluoromethane(S)	11.43	111	1201490	78.25231	ppb		0.00
Spiked Amount 25.097			Recovery = 311.795%				
38) 1,2-DCA-D4(S)	12.23	65	1017043	74.41171	ppb		0.00
Spiked Amount 24.225			Recovery = 307.167%				
56) Toluene-D8(S)	15.50	98	4141980	73.52293	ppb		0.00
Spiked Amount 25.808			Recovery = 284.881%				
64) 4-Bromofluorobenzene(S)	20.11	95	1506838	74.64914	ppb		0.00
Spiked Amount 25.459			Recovery = 293.207%				
Target Compounds							Qvalue
2) Dichlorodifluoromethane	4.07	85	884588	41.65551	ppb		100
3) Freon 114	4.33	85	557051	41.88358	ppb		95
4) Chloromethane	4.55	50	992312	37.70822	ppb		99
5) Vinyl chloride	4.81	62	582991	33.12037	ppb		100
6) 1,3-Butadiene	4.80	54	564	11.46497	ppb		93
7) Bromomethane	5.73	94	528649	41.37336	ppb		93
8) Chloroethane	5.91	64	531050	36.52439	ppb		99
9) Dichlorofluoromethane	6.01	67	1474925	36.68918	ppb		98
10) Trichlorofluoromethane	6.53	101	946797	39.68761	ppb		99
11) Acetonitrile	7.66	41	108364	171.62043	ug/l		100
12) Acrolein	7.16	56	48720	168.61606	ppb		96
13) Acetone	7.28	43	68038	41.20035	ppb	#	70
14) Freon-113	7.47	101	558655	43.47676	ppb		95
15) 1,1-DCE	7.68	96	585091	35.55608	ppb		97
16) t-Butanol	7.76	59	14424	184.59628	ppb		98
17) Methyl Acetate	8.19	43	191479	40.45406	ppb		95
18) Iodomethane	8.17	142	457316	35.89956	ppb		95
19) Acrylonitrile	8.56	53	70209	39.92472	ppb		91
20) Methylene chloride	8.48	84	561985	35.80590	ppb		98
21) Carbon disulfide	8.56	76	582016	36.40182	ppb		99
22) Methyl t-butyl ether (MtBE	8.89	73	959832	38.59078	ppb		96
23) Trans-1,2-DCE	9.10	96	690130	41.05681	ppb		94
24) Diisopropyl Ether	9.75	45	2070362	37.65279	ppb		97
25) 1,1-DCA	9.79	63	1270640	38.96588	ppb		96
26) Vinyl Acetate	9.42	43	392586	42.94098	ppb		95
27) Ethyl tert Butyl Ether	10.45	59	1446892	38.54337	ppb		98
28) MEK (2-Butanone)	10.44	43	249429	39.96063	ppb	#	93
29) Cis-1,2-DCE	10.82	96	701038	35.73563	ppb		95
30) 2,2-Dichloropropane	10.81	77	853458	36.53678	ppb		99
31) Chloroform	11.10	83	1207454	38.49020	ppb		99
32) Bromochloromethane	11.32	128	209048	38.27688	ppb		92
34) 1,1,1-TCA	11.83	97	1115691	39.10753	ppb		98
35) Cyclohexane	12.00	56	1027386	38.67839	ppb		97
36) 1,1-Dichloropropene	12.10	75	915628	37.45644	ppb		98
37) 2,2,4-Trimethylpentane	12.18	57	1588067	41.27686	ppb		96
39) Carbon Tetrachloride	12.30	117	845279	43.02829	ppb		95

(#) = qualifier out of range (m) = manual integration

Quantitation Report (Not Reviewed)

Data File : M:\CHICO\DATA\C111030\1030C22W.D Vial: 1
Acq On : 31 Oct 11 4:29 Operator: STC
Sample : Voc Std 10-30-11@40ug/L Inst : Chico
Misc : Water 10mLw/ IS:10-30-11 Multiplr: 1.00

Quant Time: Dec 2 11:19 2011 Quant Results File: CALLW.RES

Quant Method : M:\CHICO\DATA\C111030\CALLW.M (RTE Integrator)
Title : METHOD 8260
Last Update : Fri Dec 02 11:18:49 2011
Response via : Initial Calibration
DataAcq Meth : V8260

	Compound	R.T.	QIon	Response	Conc	Unit		Qvalue
40)	Tert Amyl Methyl Ether	12.34	73	1064298	37.91715	ppb		95
41)	1,2-DCA	12.37	62	609966	37.87992	ppb		96
42)	Benzene	12.50	78	2667822	37.98569	ppb		98
43)	TCE	13.53	95	750591	38.58882	ppb		94
44)	2-Pentanone	13.20	43	723061	177.71021	ppb		99
45)	1,2-Dichloropropane	13.77	63	624547	39.12345	ppb		95
46)	Bromodichloromethane	14.12	83	773755	42.42635	ppb		95
47)	Methyl Cyclohexane	13.82	83	868699	38.21751	ppb		98
48)	Dibromomethane	14.17	93	250020	39.16182	ppb		86
49)	2-Chloroethyl vinyl ether	14.57	63	173802	42.82620	ppb		95
50)	1-Bromo-2-chloroethane	14.89	63	541199	39.73256	ppb		83
51)	Cis-1,3-Dichloropropene	15.00	75	712500	40.97032	ppb		95
52)	Toluene	15.63	91	2650413	38.26138	ppb		99
53)	Trans-1,3-Dichloropropene	15.80	75	518508	41.41431	ppb		91
54)	1,1,2-TCA	16.08	83	269238	39.89943	ppb		93
57)	1,2-EDB	17.33	107	320516	42.06127	ppb		96
58)	Tetrachloroethene	16.79	164	705793	34.28742	ppb		95
59)	1-Chlorohexane	17.70	91	907243	38.27319	ppb		95
60)	1,1,1,2-Tetrachloroethane	18.16	131	576678	44.74892	ppb		93
61)	m&p-Xylene	18.35	106	2337444	76.86502	ppb		99
62)	o-Xylene	19.11	106	1156368	39.53885	ppb		94
63)	Styrene	19.13	104	1785628	40.45023	ppb		97
65)	2-Hexanone	16.10	43	145250	39.64659	ppb		96
66)	1,3-Dichloropropane	16.49	76	597192	39.74260	ppb		95
67)	Dibromochloromethane	16.97	129	435261	44.37095	ppb		90
68)	Chlorobenzene	18.10	112	1658874	38.13292	ppb		97
69)	Ethylbenzene	18.22	91	3057452	37.74292	ppb		98
70)	Bromoform	19.64	173	213787	38.87093	ppb	#	77
72)	MIBK (methyl isobutyl keto	14.67	43	228387	35.86526	ppb		86
73)	Isopropylbenzene	19.74	105	2989202	36.81944	ppb		100
74)	1,1,2,2-Tetrachloroethane	19.90	83	276570	40.69597	ppb		85
75)	1,2,3-Trichloropropane	20.15	110	27712	36.12043	ppb	#	64
76)	t-1,4-Dichloro-2-Butene	20.23	53	63970	41.51275	ppb	#	73
77)	Bromobenzene	20.47	156	676448	36.12176	ppb		94
78)	n-Propylbenzene	20.44	91	3526664	36.41330	ppb		96
79)	4-Ethyltoluene	20.64	105	2418588	36.08622	ppb		96
80)	2-Chlorotoluene	20.74	91	2324947	36.24591	ppb		98
81)	1,3,5-Trimethylbenzene	20.71	105	2435760	36.94036	ppb		97
82)	4-Chlorotoluene	20.82	91	1999529	36.20126	ppb		95
83)	Tert-Butylbenzene	21.36	119	2659556	37.25529	ppb		98
84)	1,2,4-Trimethylbenzene	21.42	105	2416954	35.09648	ppb		97
85)	Sec-Butylbenzene	21.76	105	3268087	38.17382	ppb		99
86)	p-Isopropyltoluene	21.99	119	2796609	38.13811	ppb		99
87)	Benzyl Chloride	22.42	91	391533	40.22322	ppb		98
88)	1,3-DCB	22.12	146	1411201	36.84871	ppb		99
89)	1,4-DCB	22.30	146	1322903	37.21752	ppb		96
90)	Hexachloroethane	23.60	117	488322	36.68408	ppb		93
91)	n-Butylbenzene	22.70	91	2299832	35.95845	ppb		98
92)	1,2-DCB	22.93	146	1152377	37.82968	ppb		95
93)	1,2-Dibromo-3-chloropropan	24.14	155	45893	38.08788	ppb		96
94)	1,2,4-Trichlorobenzene	25.59	180	808556	36.62946	ppb		98

(#) = qualifier out of range (m) = manual integration

Quantitation Report (Not Reviewed)

Data File : M:\CHICO\DATA\C111030\1030C22W.D Vial: 1
Acq On : 31 Oct 11 4:29 Operator: STC
Sample : Voc Std 10-30-11@40ug/L Inst : Chico
Misc : Water 10mLw/ IS:10-30-11 Multiplr: 1.00

Quant Time: Dec 2 11:19 2011 Quant Results File: CALLW.RES

Quant Method : M:\CHICO\DATA\C111030\CALLW.M (RTE Integrator)
Title : METHOD 8260
Last Update : Fri Dec 02 11:18:49 2011
Response via : Initial Calibration
DataAcq Meth : V8260

	Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
95)	Hexachlorobutadiene	25.84	223	149142	37.18851	ppb	98
96)	Naphthalene	25.94	128	1030307	37.82960	ppb	97
97)	1,2,3-Trichlorobenzene	26.29	180	633099	37.91681	ppb	96

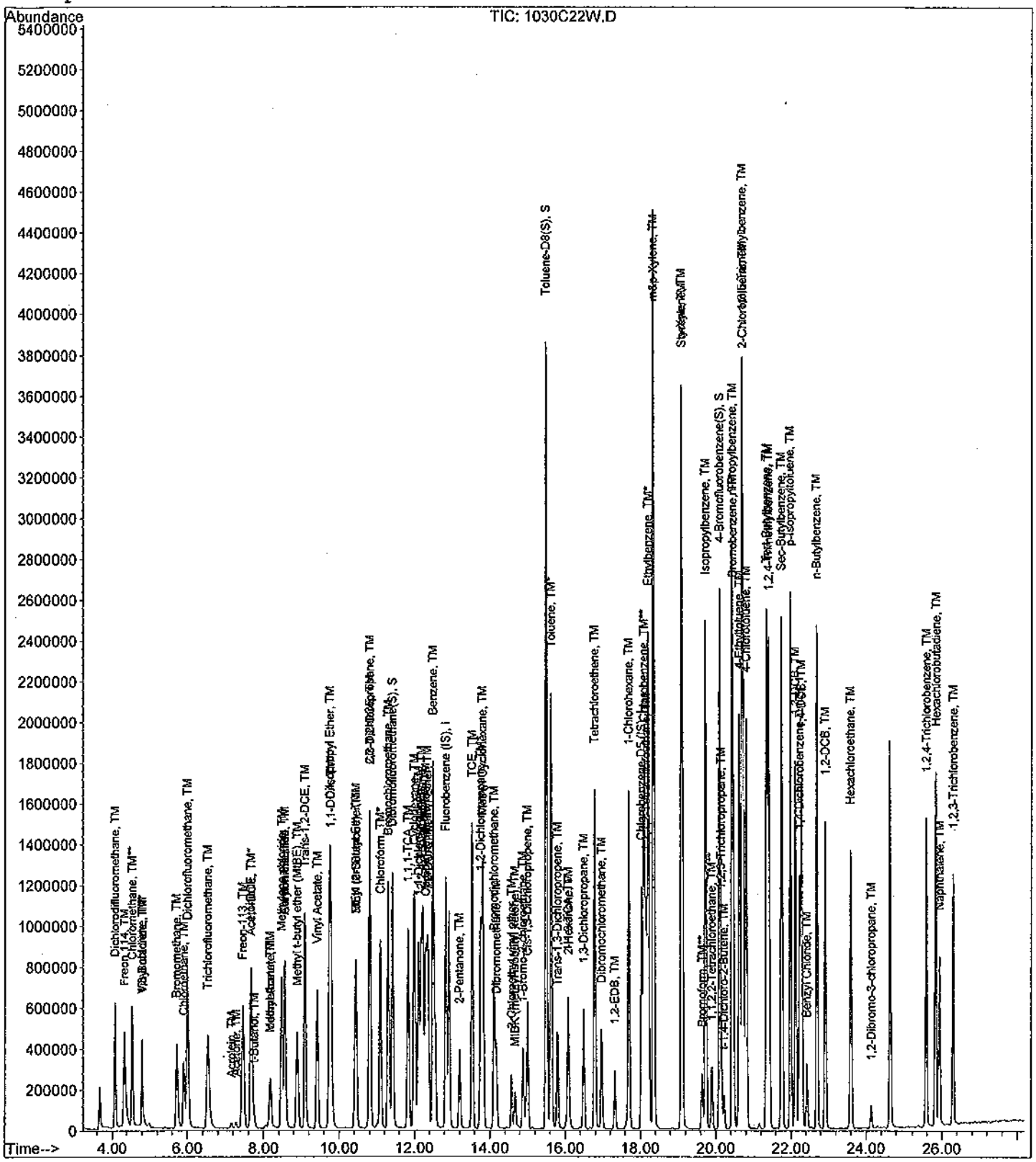
(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : M:\CHICO\DATA\C111030\1030C22W.D Vial: 1
Acq On : 31 Oct 11 4:29 Operator: STC
Sample : Voc Std 10-30-11@40ug/L Inst : Chico
Misc : Water 10mLw/ IS:10-30-11 Multiplr: 1.00

Quant Time: Dec 2 11:19 2011 Quant Results File: CALLW.RES

Method : M:\CHICO\DATA\C111030\CALLW.M (RTE Integrator)
Title : METHOD 8260
Last Update : Fri Dec 02 11:18:49 2011
Response via : Initial Calibration



Quantitation Report (Not Reviewed)

Data File : M:\CHICO\DATA\C111030\1030C23W.D Vial: 1
Acq On : 31 Oct 11 5:12 Operator: STC
Sample : Voc Std 10-30-11@100ug/L Inst : Chico
Misc : Water 10mLw/ IS:10-30-11 Multiplr: 1.00

Quant Time: Dec 2 11:19 2011 Quant Results File: CALLW.RES

Quant Method : M:\CHICO\DATA\C111030\CALLW.M (RTE Integrator)
Title : METHOD 8260
Last Update : Fri Dec 02 11:18:49 2011
Response via : Initial Calibration
DataAcq Meth : V8260

Internal Standards	R.T.	QIon	Response	Conc Units		Dev(Min)
1) Fluorobenzene (IS)	12.83	96	629184	25.00000 ppb		0.00
55) Chlorobenzene-D5 (IS)	18.04	117	438080	25.00000 ppb		0.00
71) 1,4-Dichlorobenzene-D (IS)	22.25	152	225856	25.00000 ppb		0.00
System Monitoring Compounds						
33) Dibromofluoromethane(S)	11.42	111	1562107	93.20132 ppb		0.00
Spiked Amount 25.097			Recovery = 371.359%			
38) 1,2-DCA-D4(S)	12.22	65	1315432	88.16671 ppb		0.00
Spiked Amount 24.225			Recovery = 363.946%			
56) Toluene-D8(S)	15.50	98	5498133	89.19764 ppb		0.00
Spiked Amount 25.808			Recovery = 345.617%			
64) 4-Bromofluorobenzene(S)	20.11	95	1999086	90.51343 ppb		0.00
Spiked Amount 25.459			Recovery = 355.518%			
Target Compounds						Qvalue
2) Dichlorodifluoromethane	4.07	85	2034576	87.76868 ppb		100
3) Freon 114	4.33	85	1349868	92.97677 ppb		99
4) Chloromethane	4.55	50	2507323	87.28354 ppb		98
5) Vinyl chloride	4.80	62	1338572	69.66414 ppb		97
6) 1,3-Butadiene	4.78	54	294	5.47489 ppb	#	41
7) Bromomethane	5.72	94	1399435	100.33221 ppb		96
8) Chloroethane	5.92	64	1334347	84.07184 ppb		99
9) Dichlorofluoromethane	6.00	67	3757858	85.63331 ppb		99
10) Trichlorofluoromethane	6.53	101	2271251	87.21626 ppb		100
11) Acetonitrile	7.65	41	142197	206.30447 ug/l		100
12) Acrolein	7.15	56	57928	183.65994 ppb		89
13) Acetone	7.27	43	177387	98.40233 ppb	#	80
14) Freon-113	7.46	101	1359710	98.36336 ppb		98
15) 1,1-DCE	7.69	96	1502451	83.64213 ppb		95
16) t-Butanol	7.69	59	21608	253.32966 ppb	#	70
17) Methyl Acetate	8.19	43	511503	99.99402 ppb		97
18) Iodomethane	8.16	142	1532807	102.31499 ppb		90
19) Acrylonitrile	8.57	53	190766	99.94302 ppb		88
20) Methylene chloride	8.47	84	1482941	86.55417 ppb		98
21) Carbon disulfide	8.56	76	1504256	86.18747 ppb		99
22) Methyl t-butyl ether (MtBE	8.90	73	2423605	89.26566 ppb		93
23) Trans-1,2-DCE	9.10	96	1783146	97.17967 ppb		95
24) Diisopropyl Ether	9.75	45	5216937	86.91619 ppb		95
25) 1,1-DCA	9.80	63	3188094	89.56273 ppb		96
26) Vinyl Acetate	9.42	43	966738	98.81188 ppb		94
27) Ethyl tert Butyl Ether	10.44	59	3675155	89.68576 ppb		98
28) MEK (2-Butanone)	10.42	43	621214	92.04576 ppb	#	87
29) Cis-1,2-DCE	10.82	96	1784148	83.31536 ppb		92
30) 2,2-Dichloropropane	10.81	77	2060873	80.82274 ppb		96
31) Chloroform	11.09	83	3040045	88.77564 ppb		96
32) Bromochloromethane	11.32	128	540554	90.67004 ppb		82
34) 1,1,1-TCA	11.84	97	2777762	89.19608 ppb		98
35) Cyclohexane	12.00	56	2517542	86.82519 ppb		98
36) 1,1-Dichloropropene	12.11	75	2351054	88.10577 ppb		97
37) 2,2,4-Trimethylpentane	12.18	57	4108345	99.40223 ppb		96
39) Carbon Tetrachloride	12.30	117	2194045	102.31369 ppb		94

(#) = qualifier out of range (m) = manual integration

Quantitation Report (Not Reviewed)

Data File : M:\CHICO\DATA\C111030\1030C23W.D Vial: 1
Acq On : 31 Oct 11 5:12 Operator: STC
Sample : Voc Std 10-30-11@100ug/L Inst : Chico
Misc : Water 10mLW/ IS:10-30-11 Multiplr: 1.00

Quant Time: Dec 2 11:19 2011 Quant Results File: CALLW.RES

Quant Method : M:\CHICO\DATA\C111030\CALLW.M (RTE Integrator)
Title : METHOD 8260
Last Update : Fri Dec 02 11:18:49 2011
Response via : Initial Calibration
DataAcq Meth : V8260

	Compound	R.T.	QIon	Response	Conc Unit		Qvalue
40)	Tert Amyl Methyl Ether	12.35	73	2816735	91.92902 ppb		99
41)	1,2-DCA	12.38	62	1561150	88.81419 ppb		98
42)	Benzene	12.49	78	6991851	91.19889 ppb		98
43)	TCE	13.54	95	1880975	88.58811 ppb		90
44)	2-Pentanone	13.20	43	881325	198.43021 ppb		99
45)	1,2-Dichloropropane	13.76	63	1582643	90.82163 ppb		96
46)	Bromodichloromethane	14.11	83	1954559	98.17824 ppb	#	91
47)	Methyl Cyclohexane	13.82	83	2215169	89.27589 ppb		99
48)	Dibromomethane	14.17	93	660037	94.70886 ppb		92
49)	2-Chloroethyl vinyl ether	14.57	63	484281	109.31668 ppb		95
50)	1-Bromo-2-chloroethane	14.88	63	1381119	92.88701 ppb		92
51)	Cis-1,3-Dichloropropene	15.00	75	1816087	95.66550 ppb		93
52)	Toluene	15.64	91	6789315	89.78572 ppb		99
53)	Trans-1,3-Dichloropropene	15.80	75	1379809	100.95973 ppb		94
54)	1,1,2-TCA	16.08	83	670730	91.05678 ppb		90
57)	1,2-EDB	17.32	107	832657	99.86703 ppb		98
58)	Tetrachloroethene	16.79	164	1730703	76.84281 ppb		95
59)	1-Chlorohexane	17.71	91	2353794	90.75337 ppb		91
60)	1,1,1,2-Tetrachloroethane	18.16	131	1508942	107.01509 ppb		94
61)	m&p-Xylene	18.36	106	6180125	185.74112 ppb		97
62)	o-Xylene	19.10	106	3002956	93.84264 ppb		92
63)	Styrene	19.12	104	4577224	94.76663 ppb		100
65)	2-Hexanone	16.10	43	375289	93.62219 ppb		94
66)	1,3-Dichloropropane	16.49	76	1470502	89.43988 ppb		98
67)	Dibromochloromethane	16.97	129	1191759	111.03520 ppb		92
68)	Chlorobenzene	18.11	112	4271113	89.73276 ppb		97
69)	Ethylbenzene	18.22	91	8013287	90.40861 ppb		94
70)	Bromoform	19.64	173	616423	100.58038 ppb	#	81
72)	MIBK (methyl isobutyl keto	14.67	43	581084	90.50195 ppb		93
73)	Isopropylbenzene	19.74	105	7614687	93.02299 ppb		99
74)	1,1,2,2-Tetrachloroethane	19.90	83	720253	105.11090 ppb		88
75)	1,2,3-Trichloropropane	20.16	110	78648	101.50443 ppb	#	68
76)	t-1,4-Dichloro-2-Butene	20.22	53	179966	115.82759 ppb	#	75
77)	Bromobenzene	20.48	156	1766849	93.57294 ppb		97
78)	n-Propylbenzene	20.45	91	9038917	92.56114 ppb		95
79)	4-Ethyltoluene	20.64	105	6293560	93.13058 ppb		93
80)	2-Chlorotoluene	20.74	91	5791730	89.55104 ppb		98
81)	1,3,5-Trimethylbenzene	20.72	105	6243237	93.90589 ppb		96
82)	4-Chlorotoluene	20.81	91	5272321	94.67040 ppb		97
83)	Tert-Butylbenzene	21.36	119	6787695	94.30127 ppb		99
84)	1,2,4-Trimethylbenzene	21.41	105	6234859	89.79211 ppb		97
85)	Sec-Butylbenzene	21.76	105	8294079	96.08521 ppb		99
86)	p-Isopropyltoluene	21.99	119	7113742	96.21480 ppb		98
87)	Benzyl Chloride	22.43	91	1035237	105.47867 ppb		97
88)	1,3-DCB	22.13	146	3600903	93.25266 ppb		96
89)	1,4-DCB	22.29	146	3434577	95.83166 ppb		96
90)	Hexachloroethane	23.60	117	1413252	101.62319 ppb		97
91)	n-Butylbenzene	22.70	91	5960168	92.42293 ppb		99
92)	1,2-DCB	22.92	146	2973338	96.80520 ppb		95
93)	1,2-Dibromo-3-chloropropan	24.14	155	125399	101.01251 ppb		95
94)	1,2,4-Trichlorobenzene	25.58	180	2057289	92.43409 ppb		98

(#) = qualifier out of range (m) = manual integration

Quantitation Report (Not Reviewed)

Data File : M:\CHICO\DATA\C111030\1030C23W.D Vial: 1
Acq On : 31 Oct 11 5:12 Operator: STC
Sample : Voc Std 10-30-11@100ug/L Inst : Chico
Misc : Water 10mLw/ IS:10-30-11 Multiplr: 1.00

Quant Time: Dec 2 11:19 2011 Quant Results File: CALLW.RES

Quant Method : M:\CHICO\DATA\C111030\CALLW.M (RTE Integrator)
Title : METHOD 8260
Last Update : Fri Dec 02 11:18:49 2011
Response via : Initial Calibration
DataAcq Meth : V8260

Compound	R.T.	QIon	Response	Conc Unit	Qvalue
95) Hexachlorobutadiene	25.83	223	370844	91.70996 ppb	98
96) Naphthalene	25.93	128	2644717	96.30764 ppb	98
97) 1,2,3-Trichlorobenzene	26.30	180	1614041	95.87185 ppb	100

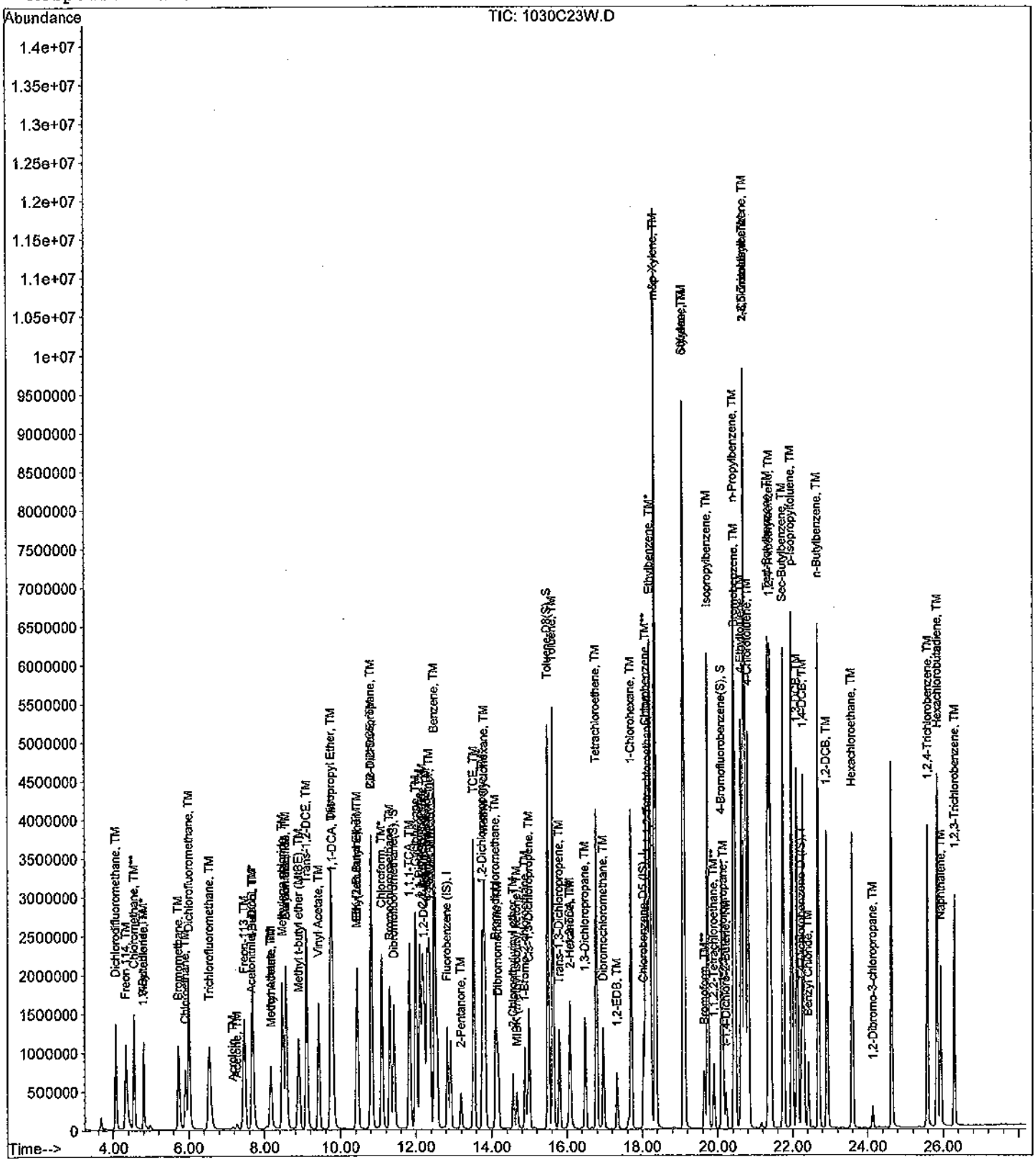
(#) = qualifier out of range (m) = manual integration

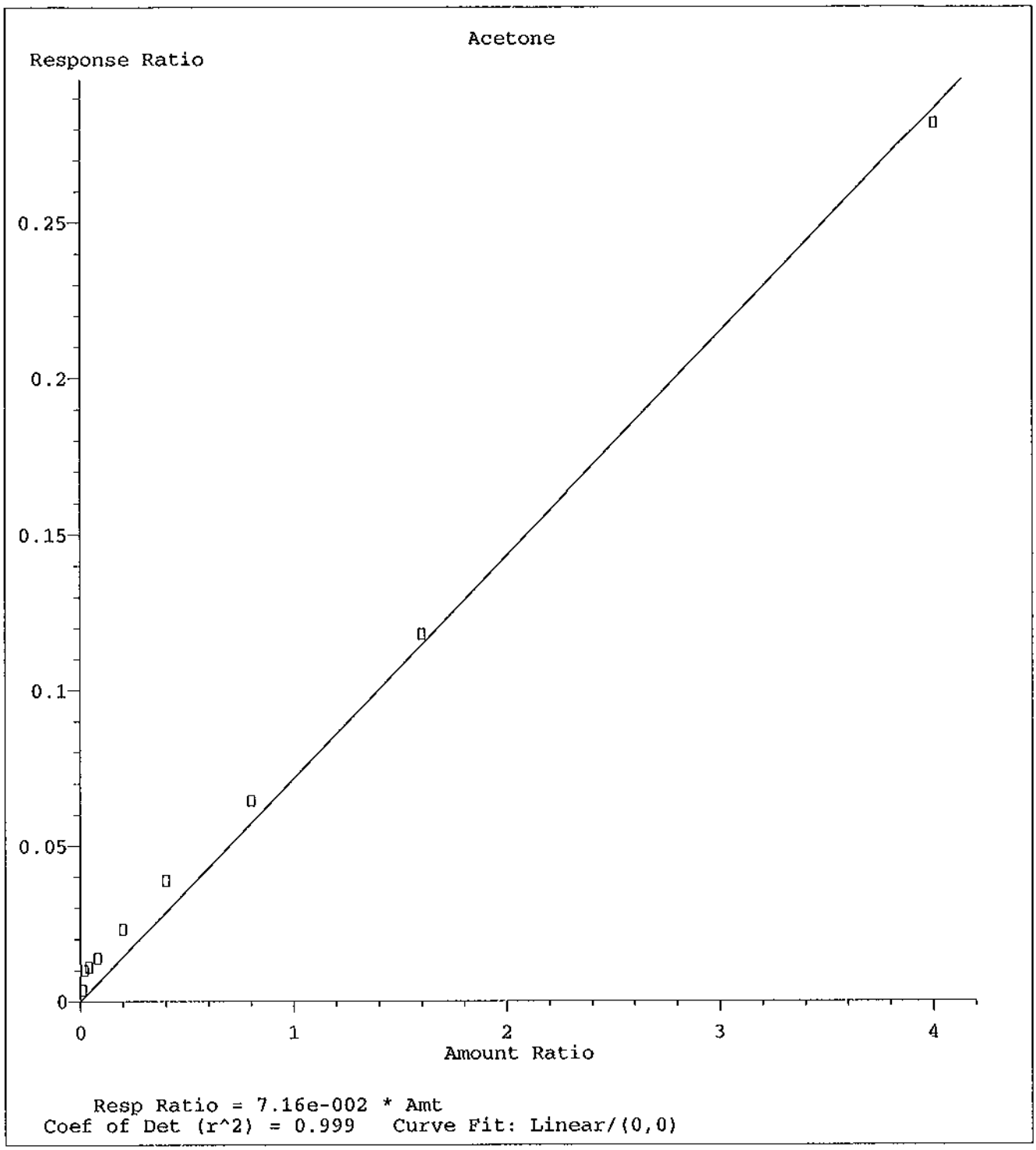
Quantitation Report

Data File : M:\CHICO\DATA\C111030\1030C23W.D Vial: 1
Acq On : 31 Oct 11 5:12 Operator: STC
Sample : Voc Std 10-30-11@100ug/L Inst : Chico
Misc : Water 10mLw/ IS:10-30-11 Multiplr: 1.00

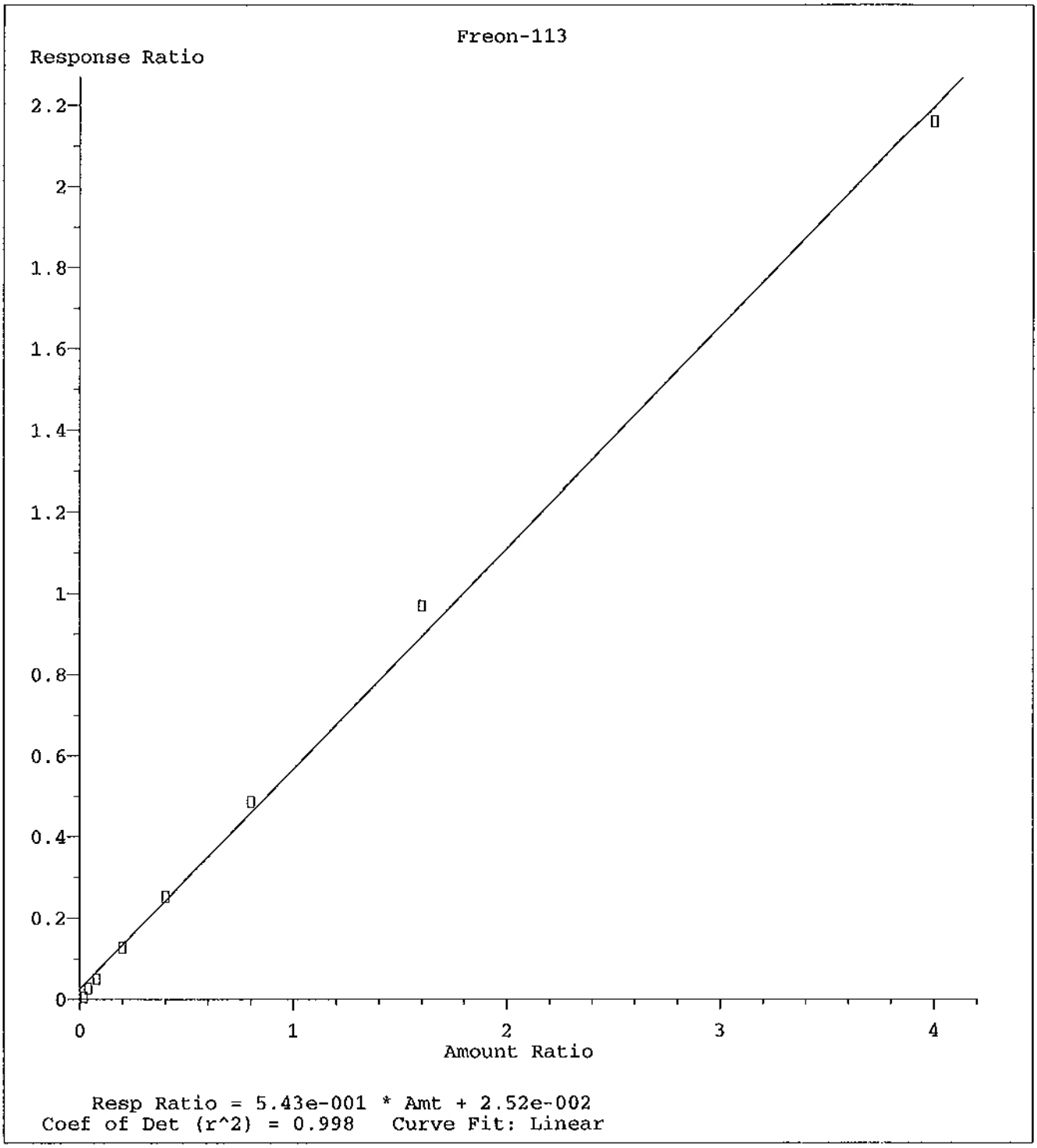
Quant Time: Dec 2 11:19 2011 Quant Results File: CALLW.RES

Method : M:\CHICO\DATA\C111030\CALLW.M (RTE Integrator)
Title : METHOD 8260
Last Update : Fri Dec 02 11:18:49 2011
Response via : Initial Calibration

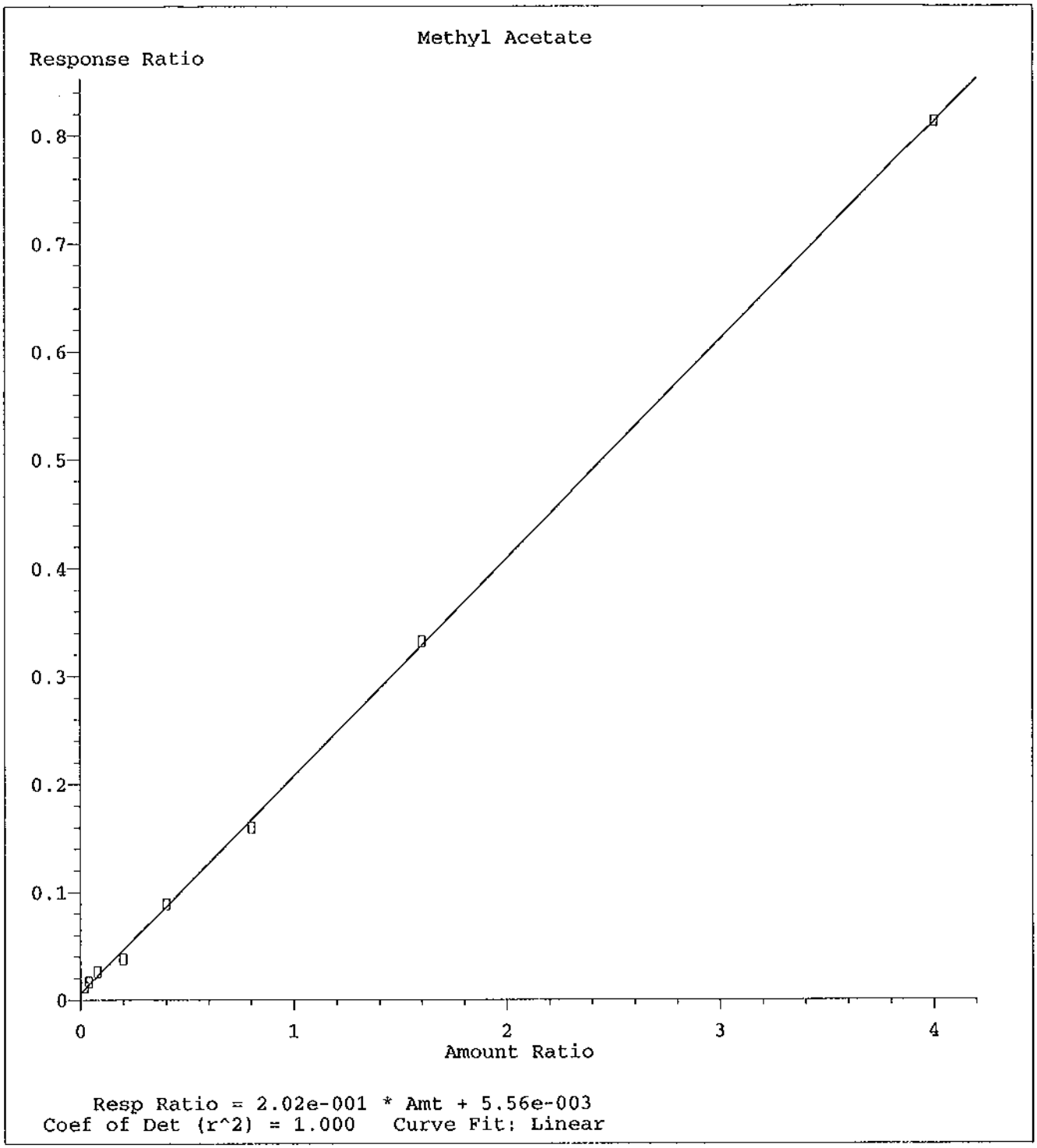




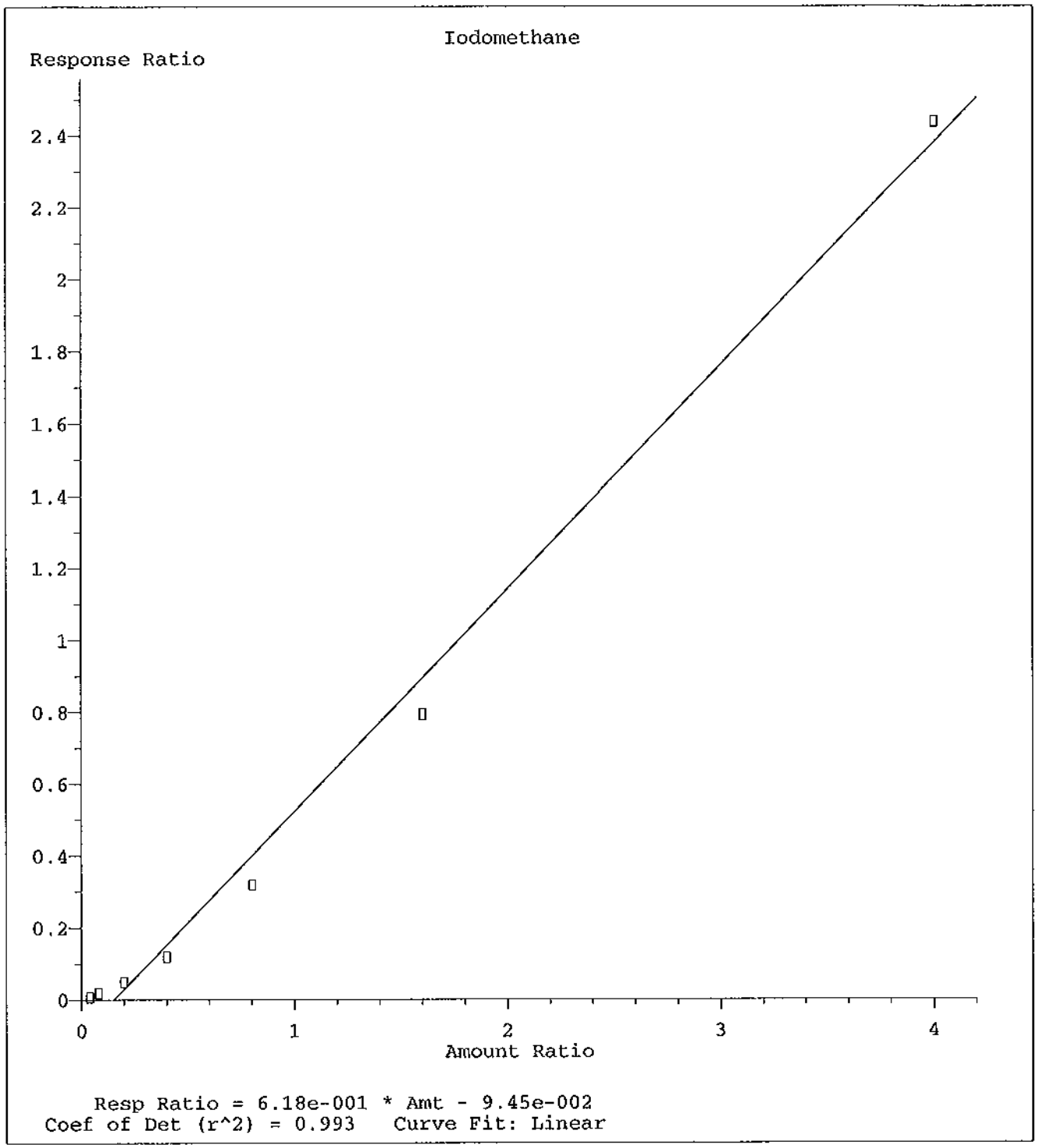
Method Name: M:\CHICO\DATA\C111030\CALLW.M
Calibration Table Last Updated: Fri Dec 02 11:32:50 2011



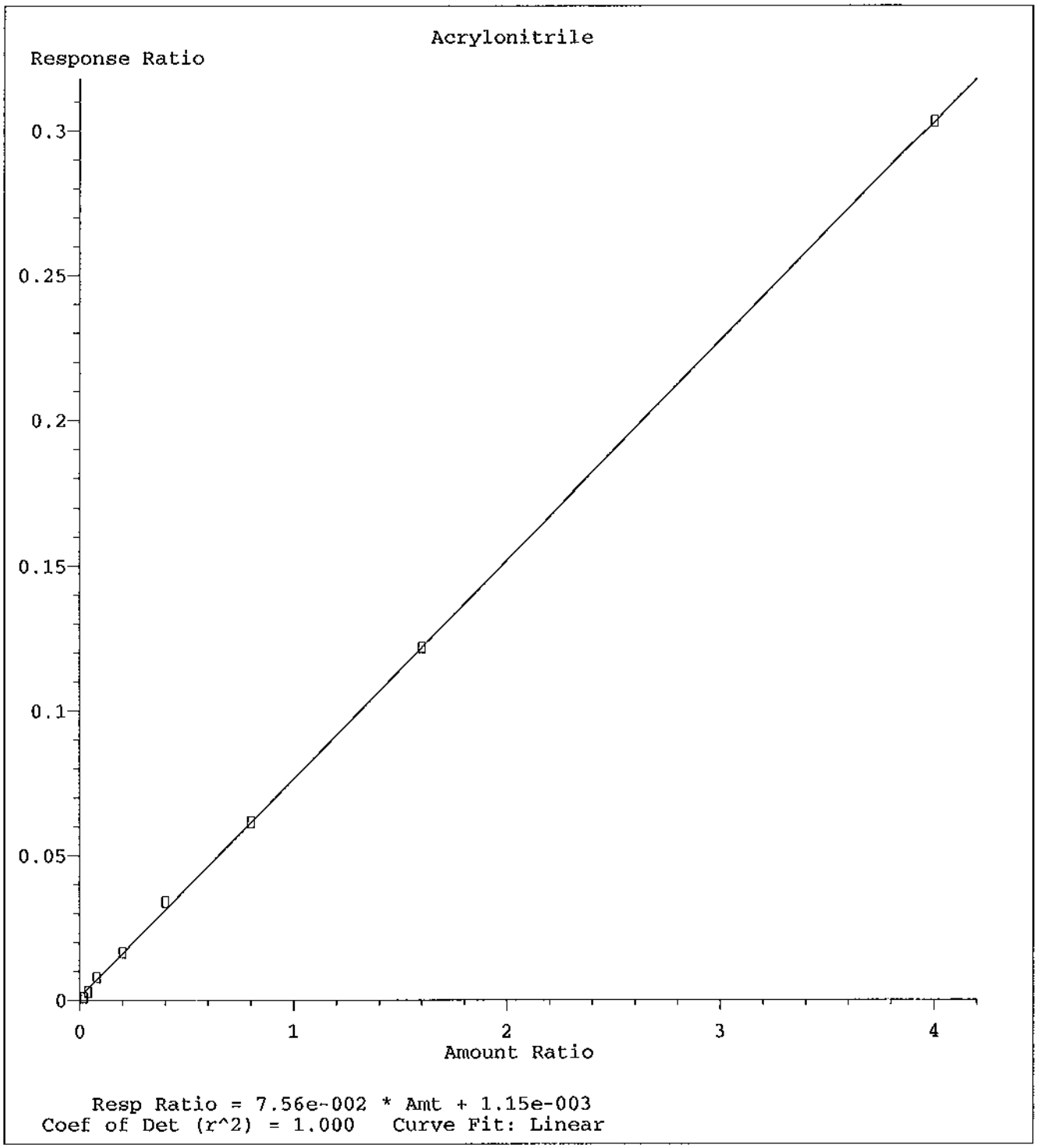
Method Name: M:\CHICO\DATA\C111030\CALLW.M
Calibration Table Last Updated: Fri Dec 02 11:32:50 2011



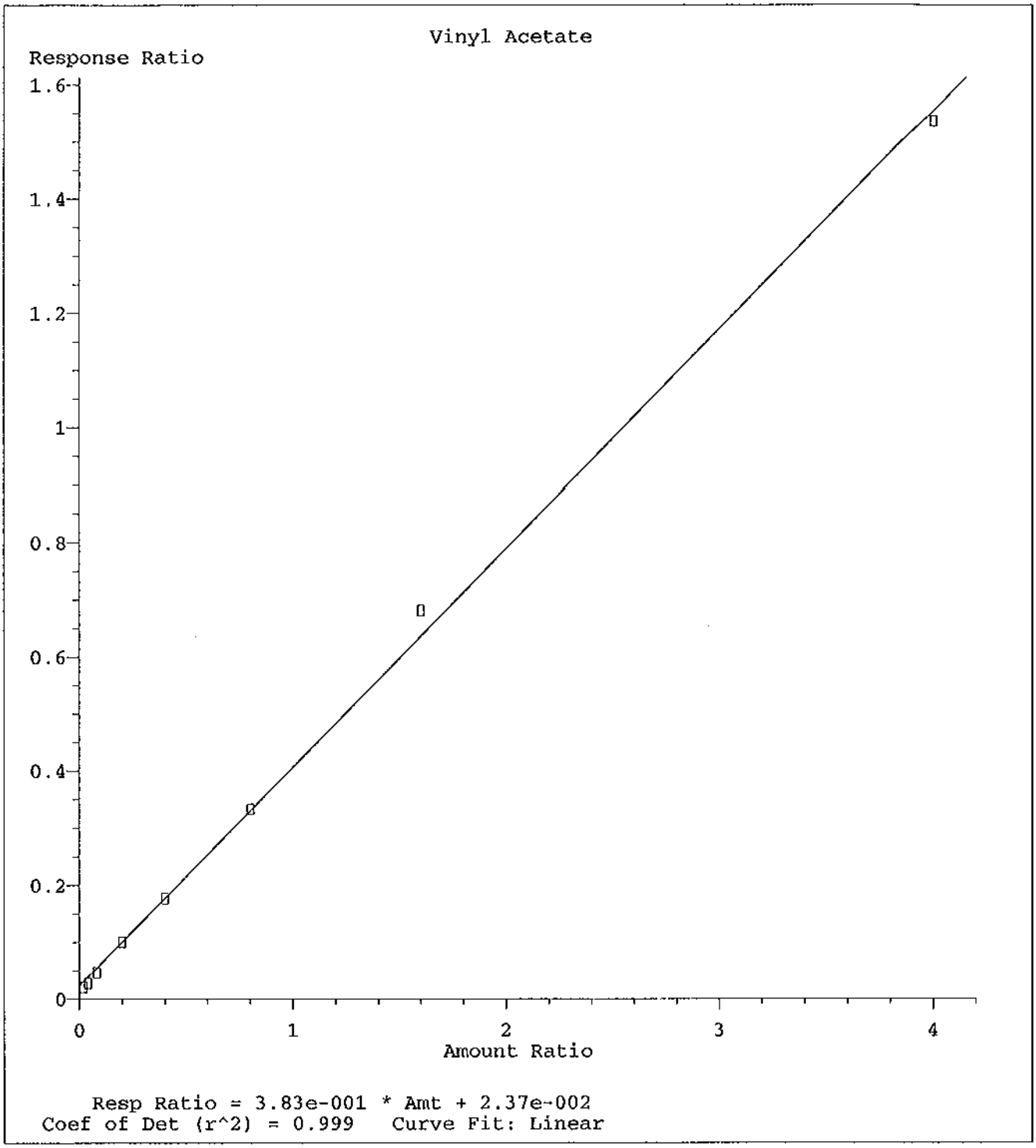
Method Name: M:\CHICO\DATA\C111030\CALLW.M
Calibration Table Last Updated: Fri Dec 02 11:32:50 2011



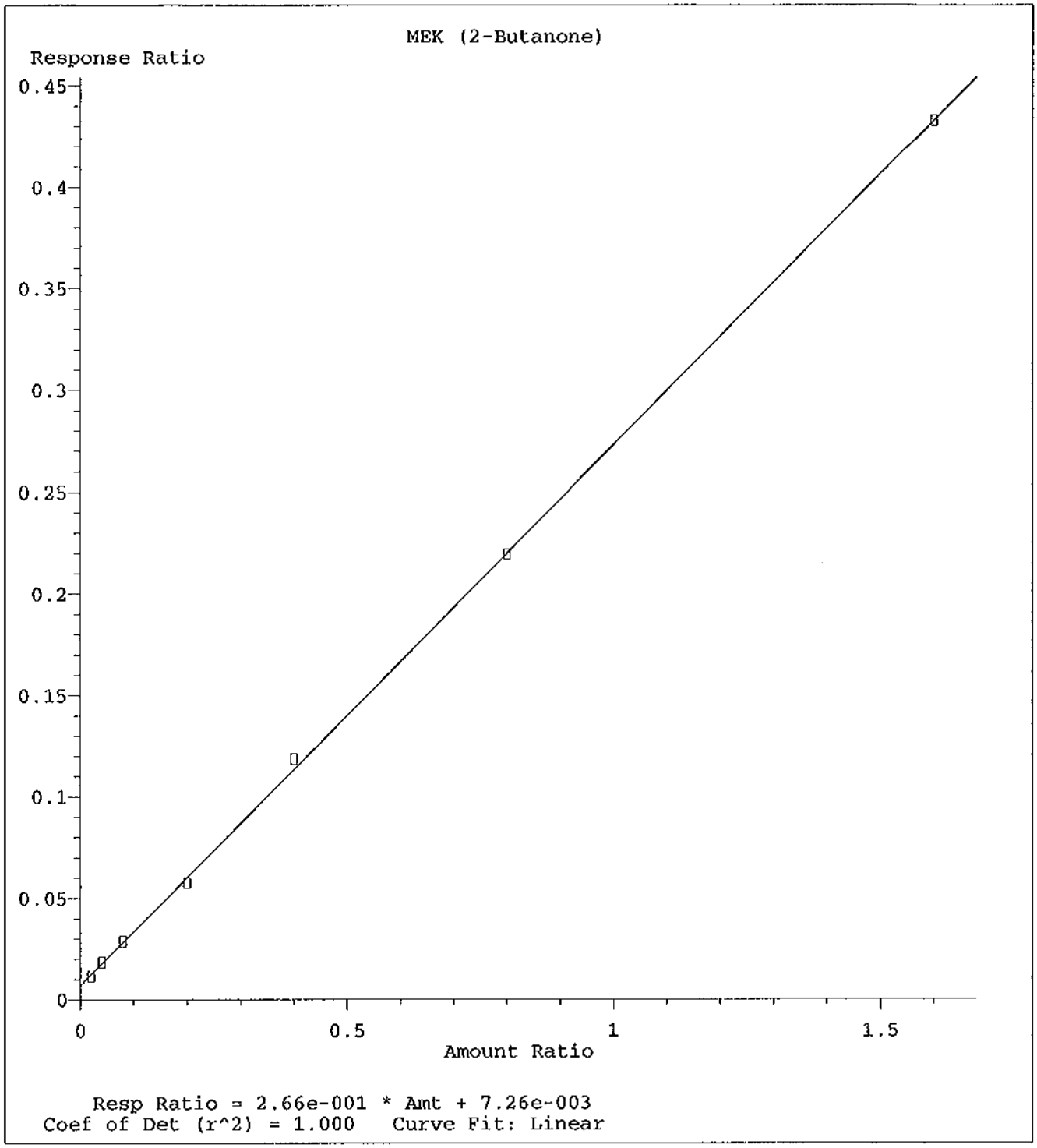
Method Name: M:\CHICO\DATA\C111030\CALLW.M
Calibration Table Last Updated: Fri Dec 02 11:32:50 2011



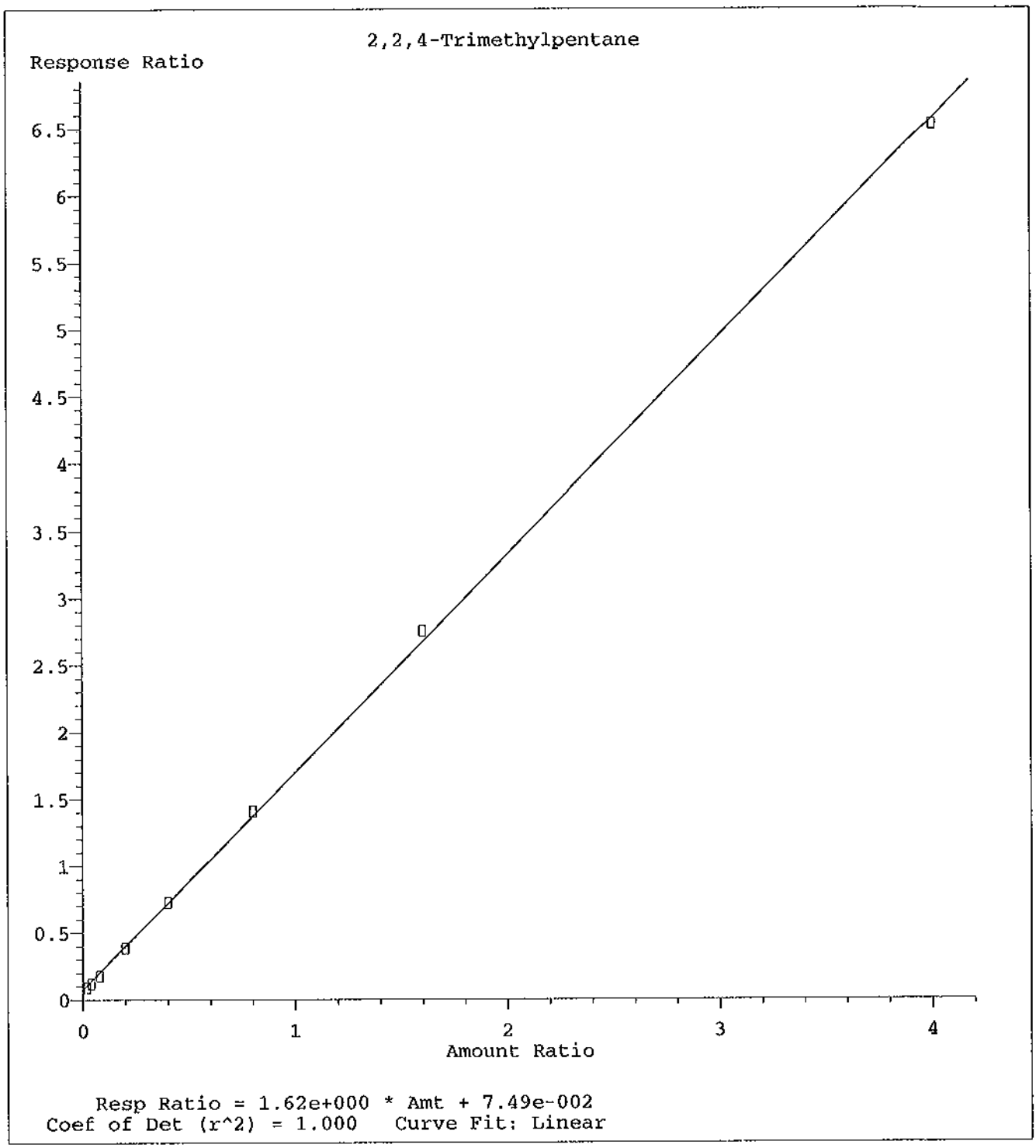
Method Name: M:\CHICO\DATA\C111030\CALLW.M
Calibration Table Last Updated: Fri Dec 02 11:32:50 2011



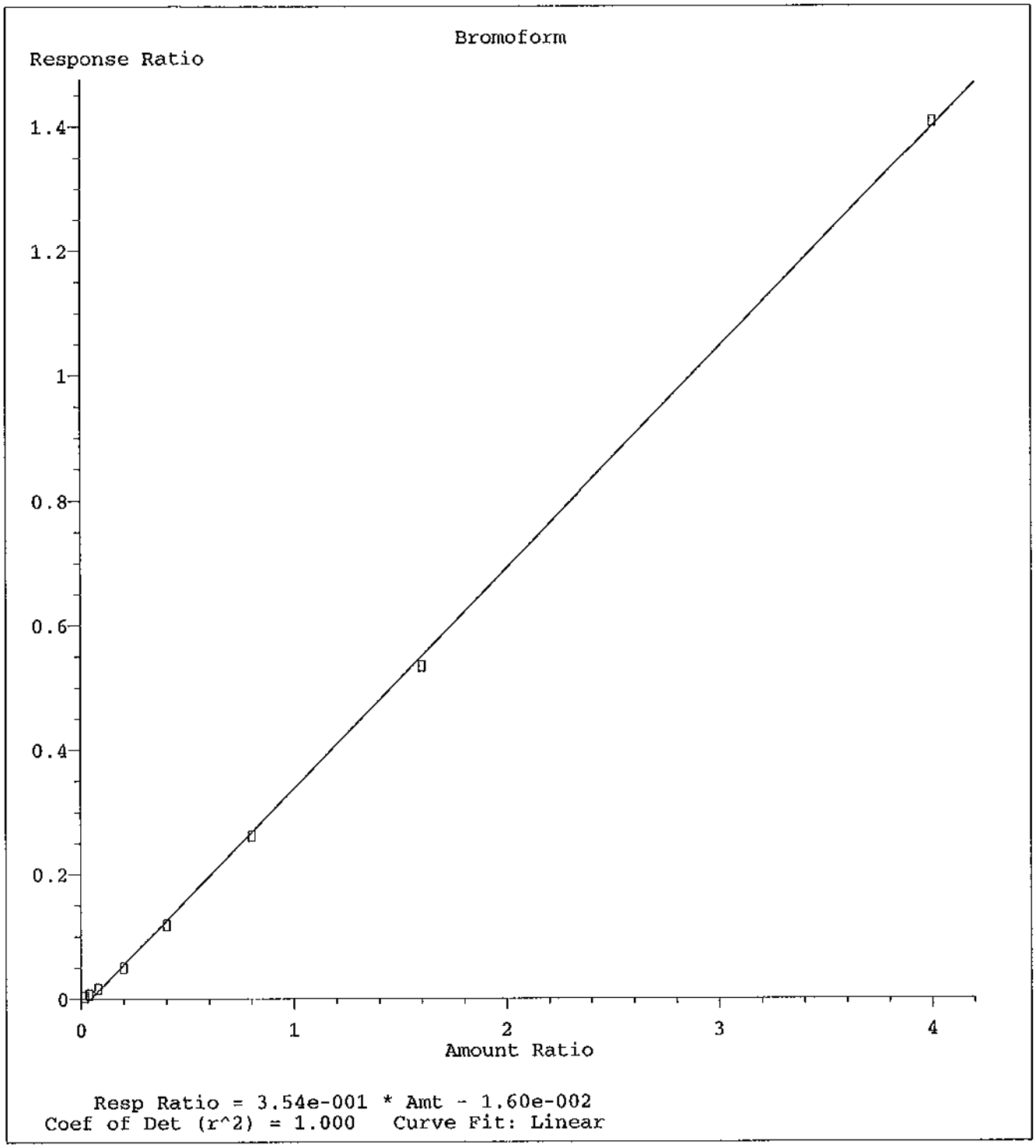
Method Name: M:\CHICO\DATA\C111030\CALLW.M
Calibration Table Last Updated: Fri Dec 02 11:32:50 2011



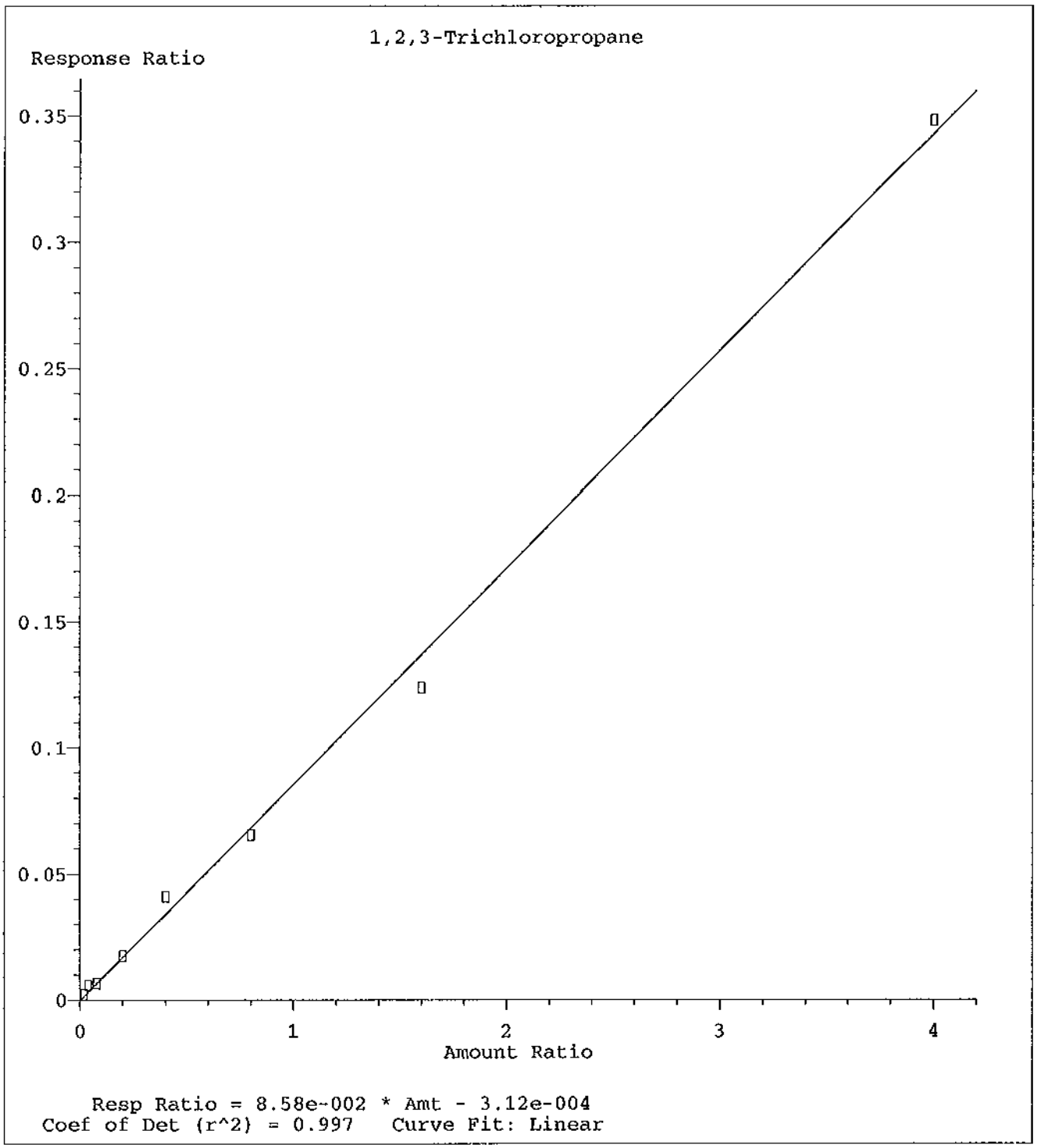
Method Name: M:\CHICO\DATA\C111030\CALLW.M
Calibration Table Last Updated: Fri Dec 02 11:32:50 2011



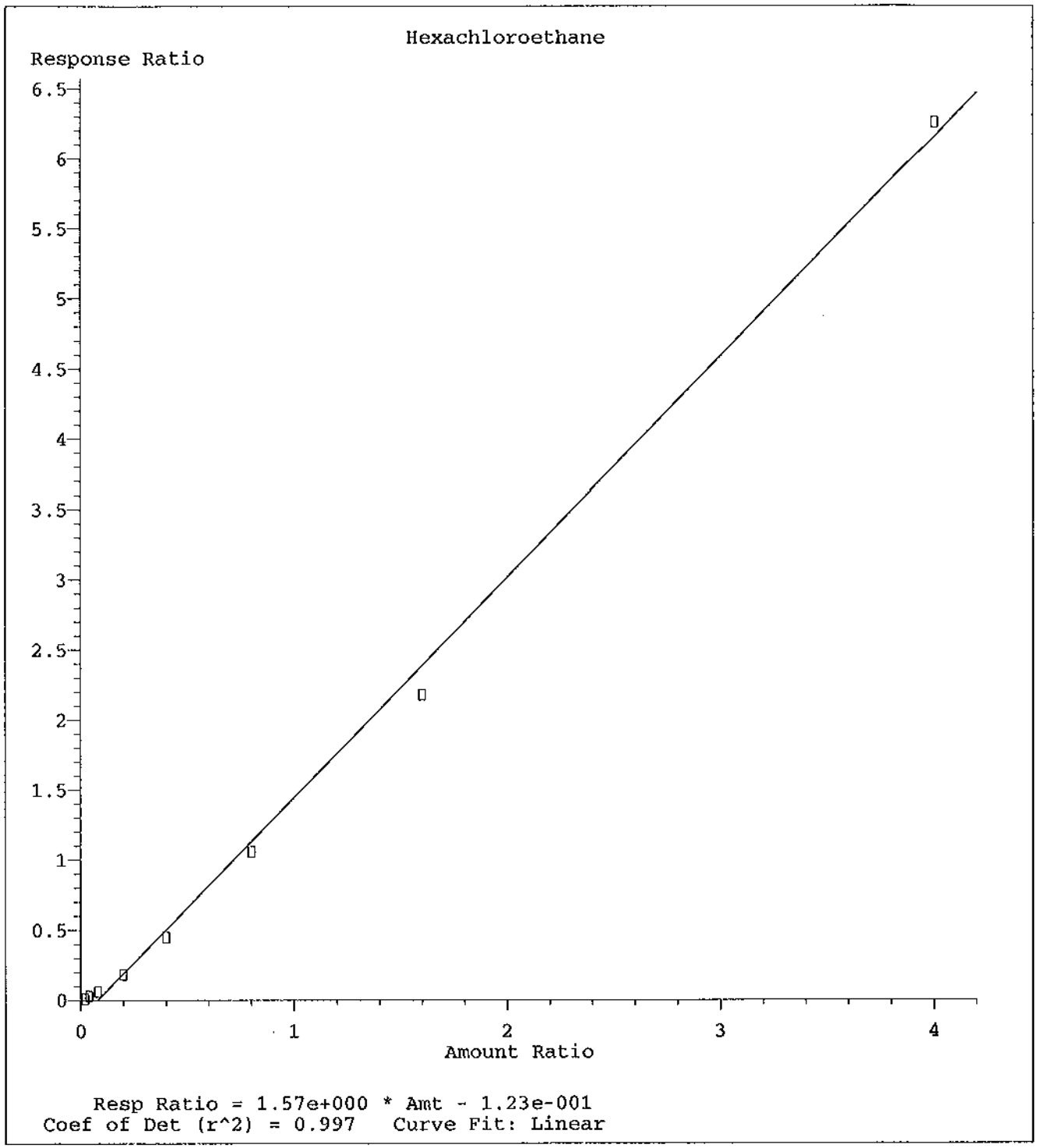
Method Name: M:\CHICO\DATA\C111030\CALLW.M
Calibration Table Last Updated: Fri Dec 02 11:32:50 2011



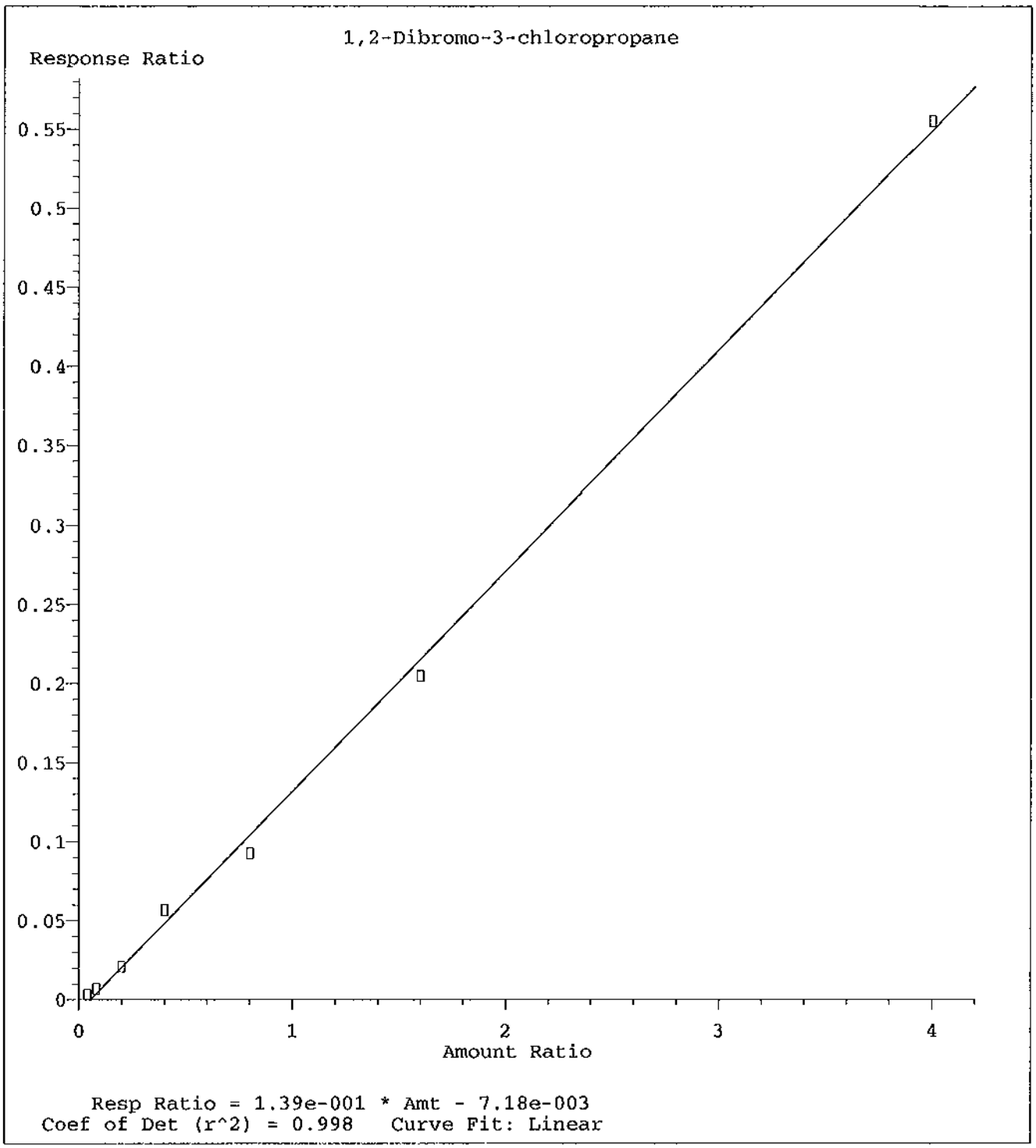
Method Name: M:\CHICO\DATA\C111030\CALLW.M
Calibration Table Last Updated: Fri Dec 02 11:32:50 2011



Method Name: M:\CHICO\DATA\C111030\CALLW.M
Calibration Table Last Updated: Fri Dec 02 11:32:50 2011



Method Name: M:\CHICO\DATA\C111030\CALLW.M
Calibration Table Last Updated: Fri Dec 02 11:32:50 2011



Method Name: M:\CHICO\DATA\C111030\CALLW.M
Calibration Table Last Updated: Fri Dec 02 11:32:50 2011

VOLATILE ORGANIC ANALYSIS BY
EPA METHOD 8260B

Form 7

Second Source Calibration

Lab Name: APPL, Inc.
Case No:
Matrix:

SDG No:
Date Analyzed: 10/31/11
Instrument: Chico
Initial Cal. Date: 10/30/11
Data File: 1030C28W.D

		Compound	MEAN	CCRF	%D		%Drift	
1	I	Fluorobenzene (IS)	ISTD			I		
2	TM	Dichlorodifluoromethane	0.9211	0.9249	0.42	TM		
3	TM	Freon 114	0.5769	0.6094	5.6	TM		
4	TM**	Chloromethane	1.141	1.076	5.8	TM**		
5	TM*	Vinyl chloride	0.7635	0.7773	1.8	TM*		
6	TML	1,3-Butadiene	0.0000	0.0015	0.00	TML		
7	TM	Bromomethane	0.5542	0.4961	10	TM		
8	TM	Chloroethane	0.6306	0.5609	11	TM		
9	TM	Dichlorofluoromethane	1.744	1.607	7.8	TM		
10	TM	Trichlorofluoromethane	1.035	1.006	2.8	TM		
11		Acetonitrile	0.0274	0.0263	4.1			
12	TM	Acrolein	0.0125	0.0115	8.4	TM		
13	TML	Acetone	0.1859	0.0931	50	TML	30	*
14	TML	Freon-113	0.5715	0.5758	0.75	TML	5.5	
15	TM*	1,1-DCE	0.7137	0.6303	12	TM*		
16	TM	t-Butanol	0.0034	0.0035	3.5	TM		
17	TML	Methyl Acetate	0.2927	0.2030	31	TML	6.3	
18	TML	Iodomethane	0.3500	0.4253	21	TML	7.0	
19	TML	Acrylonitrile	0.0764	0.0746	2.3	TML	5.1	
20	TM	Methylene chloride	0.6808	0.6159	9.5	TM		
21	TM	Carbon disulfide	0.6935	0.6450	7.0	TM		
22	TM	Methyl t-butyl ether (MtBE)	1.079	1.046	3.1	TM		
23	TM	Trans-1,2-DCE	0.8280	0.7496	9.5	TM		
24	TM	Diisopropyl Ether	2.385	2.306	3.3	TM		
25	TM**	1,1-DCA	1.414	1.411	0.23	TM**		
26	TML	Vinyl Acetate	0.5623	0.4364	22	TML	1.5	
27	TM	Ethyl tert Butyl Ether	1.628	1.646	1.1	TM		
28	TML	MEK (2-Butanone)	0.3591	0.2972	17	TML	4.8	
29	TM	Cis-1,2-DCE	0.8509	0.7812	8.2	TM		
30	TM	2,2-Dichloropropane	1.013	0.8669	14	TM		
31	TM*	Chloroform	1.361	1.332	2.1	TM*		
32	TM	Bromochloromethane	0.2369	0.2434	2.8	TM		
33	S	Dibromofluoromethane(S)	0.6660	0.6700	0.60	S		
34	TM	1,1,1-TCA	1.237	1.182	4.5	TM		
35	TM	Cyclohexane	1.152	1.120	2.8	TM		
36	TM	1,1-Dichloropropene	1.060	0.9998	5.7	TM		
37	TML	2,2,4-Trimethylpentane	2.316	1.725	26	TML	5.3	
38	S	1,2-DCA-D4(S)	0.5928	0.5784	2.4	S		
39	TM	Carbon Tetrachloride	0.8521	0.8321	2.3	TM		
40	TM	Tert Amyl Methyl Ether	1.217	1.216	0.15	TM		
		Average			8.3			

VOLATILE ORGANIC ANALYSIS BY
EPA METHOD 8260B

Form 7

Second Source Calibration

Lab Name: APPL, Inc.
Case No:
Matrix: 0

SDG No:
Date Analyzed: 10/31/11
Instrument: Chico
Cal. Date: 10/30/11
Data File: 1030C28W.D

		Compound	MEAN	CCRF	%D		%Drift
41	TM	1,2-DCA	0.6984	0.6709	3.9	TM	
42	TM	Benzene	3.046	2.887	5.2	TM	
43	TM	TCE	0.8437	0.8134	3.6	TM	
44	TM	2-Pentanone	0.1765	0.1769	0.25	TM	
45	TM*	1,2-Dichloropropane	0.6924	0.6963	0.57	TM*	
46	TM	Bromodichloromethane	0.7910	0.7910	0.01	TM	
47	TM	Methyl Cyclohexane	0.9859	0.9347	5.2	TM	
48	TM	Dibromomethane	0.2769	0.2811	1.5	TM	
49	TM	2-Chloroethyl vinyl ether	0.1760	0.1764	0.23	TM	
50	TM	1-Bromo-2-chloroethane	0.5908	0.6129	3.7	TM	
51	TM	Cis-1,3-Dichloropropene	0.7543	0.7571	0.37	TM	
52	TM*	Toluene	3.005	2.824	6.0	TM*	
53	TM	Trans-1,3-Dichloropropene	0.5430	0.5270	3.0	TM	
54	TM	1,1,2-TCA	0.2927	0.2901	0.90	TM	
55	I	Chlorobenzene-D5 (IS)	ISTD			I	
56	S	Toluene-D8(S)	3.518	3.782	7.5	S	
57	TM	1,2-EDB	0.4758	0.5074	6.6	TM	
58	TM	Tetrachloroethene	1.285	1.296	0.81	TM	
59	TM	1-Chlorohexane	1.480	1.493	0.86	TM	
60	TM	1,1,1,2-Tetrachloroethane	0.8047	0.8891	10	TM	
61	TM	m&p-Xylene	1.899	1.889	0.49	TM	
62	TM	o-Xylene	1.826	1.894	3.7	TM	
63	TM	Styrene	2.756	2.935	6.5	TM	
64	S	4-Bromofluorobenzene(S)	1.260	1.394	11	S	
65	TM	2-Hexanone	0.2288	0.2487	8.7	TM	
66	TM	1,3-Dichloropropane	0.9383	0.9463	0.86	TM	
67	TM	Dibromochloromethane	0.6125	0.6760	10	TM	
68	TM**	Chlorobenzene	2.716	2.809	3.4	TM**	
69	TM*	Ethylbenzene	5.058	5.094	0.71	TM*	
70	TM**L	Bromoform	0.2607	0.2895	11	TM**L	6.8
71	I	1,4-Dichlorobenzene-D (IS)	ISTD			I	
72	TM	MIBK (methyl isobutyl ketone)	0.7107	0.6949	2.2	TM	
73	TM	Isopropylbenzene	9.061	8.888	1.9	TM	
74	TM**	1,1,2,2-Tetrachloroethane	0.7585	0.7787	2.7	TM**	
75	TML	1,2,3-Trichloropropane	0.0967	0.1006	4.1	TML	18
76	TM	t-1,4-Dichloro-2-Butene	0.1720	0.1758	2.2	TM	
77	TM	Bromobenzene	2.090	2.104	0.64	TM	
78	TM	n-Propylbenzene	10.8	10.6	2.1	TM	
79	TM	4-Ethyltoluene	7.480	6.973	6.8	TM	
80	TM	2-Chlorotoluene	7.159	7.027	1.8	TM	
		Average			3.7		

VOLATILE ORGANIC ANALYSIS BY
EPA METHOD 8260B

Form 7

Second Source Calibration

Lab Name: APPL, Inc.
Case No:
Matrix: 0

SDG No:
Date Analyzed: 10/31/11
Instrument: Chico
Cal. Date: 10/30/11
Data File: 1030C28W.D

		Compound	MEAN	CCRF	%D		%Drift
81	TM	1,3,5-Trimethylbenzene	7.359	7.258	1.4	TM	
82	TM	4-Chlorotoluene	6.164	5.899	4.3	TM	
83	TM	Tert-Butylbenzene	7.967	7.897	0.88	TM	
84	TM	1,2,4-Trimethylbenzene	7.686	7.097	7.7	TM	
85	TM	Sec-Butylbenzene	9.555	9.679	1.3	TM	
86	TM	p-Isopropyltoluene	8.184	8.031	1.9	TM	
87	TM	Benzyl Chloride	1.086	0.9559	12	TM	
88	TM	1,3-DCB	4.274	4.133	3.3	TM	
89	TM	1,4-DCB	3.967	3.945	0.56	TM	
90	TML	Hexachloroethane	1.021	1.085	6.2	TML	11
91	TM	n-Butylbenzene	7.138	6.849	4.1	TM	
92	TM	1,2-DCB	3.400	3.438	1.1	TM	
93	TML	1,2-Dibromo-3-chloropropane	0.1148	0.1240	8.0	TML	2.0
94	TM	1,2,4-Trichlorobenzene	2.464	2.504	1.7	TM	
95	TM	Hexachlorobutadiene	0.4476	0.4720	5.5	TM	
96	TM	Naphthalene	3.040	3.207	5.5	TM	
97	TM	1,2,3-Trichlorobenzene	1.864	2.025	8.7	TM	
98							
99							
100							
101							
102							
103							
104							
105							
106							
107							
108							
109							
110							
111							
112							
113							
114							
115							
116							
117							
118							
119							
120							
		Average			4.4		

Quantitation Report (Not Reviewed)

Data File : M:\CHICO\DATA\C111030\1030C28W.D Vial: 1
Acq On : 31 Oct 11 8:48 Operator: STC
Sample : 111030A LCS-1WC (SS) Inst : Chico
Misc : Water 10mLw/ IS&S:10-30/10-26-11 Multiplr: 1.00

Quant Time: Dec 2 11:35 2011 Quant Results File: CALLW.RES

Quant Method : M:\CHICO\DATA\C111030\CALLW.M (RTE Integrator)
Title : METHOD 8260
Last Update : Fri Dec 02 11:32:50 2011
Response via : Initial Calibration
DataAcq Meth : V8260

Internal Standards	R.T.	QIon	Response	Conc Units		Dev(Min)
1) Fluorobenzene (IS)	12.84	96	600576	25.00000 ppb		0.00
55) Chlorobenzene-D5 (IS)	18.04	117	389760	25.00000 ppb		0.00
71) 1,4-Dichlorobenzene-D (IS)	22.24	152	212800	25.00000 ppb		0.00
System Monitoring Compounds						
33) Dibromofluoromethane(S)	11.42	111	402364	25.15012 ppb		0.00
Spiked Amount 25.097			Recovery	= 100.210%		
38) 1,2-DCA-D4(S)	12.23	65	347346	24.38980 ppb		0.00
Spiked Amount 24.225			Recovery	= 100.680%		
56) Toluene-D8(S)	15.50	98	1474138	26.88019 ppb		0.00
Spiked Amount 25.808			Recovery	= 104.153%		
64) 4-Bromofluorobenzene(S)	20.11	95	543410	27.65447 ppb		0.00
Spiked Amount 25.459			Recovery	= 108.620%		
Target Compounds						Qvalue
2) Dichlorodifluoromethane	4.07	85	222193	10.04166 ppb		100
3) Freon 114	4.33	85	146402	10.56428 ppb		93
4) Chloromethane	4.55	50	258385	9.42321 ppb		99
5) Vinyl chloride	4.82	62	186738	10.18146 ppb		96
7) Bromomethane	5.72	94	119189	8.95228 ppb		94
8) Chloroethane	5.91	64	134747	8.89428 ppb		97
9) Dichlorofluoromethane	6.01	67	386134	9.21828 ppb		97
10) Trichlorofluoromethane	6.52	101	241622	9.72027 ppb		99
11) Acetonitrile	7.64	41	78885	119.90087 ug/l		100
12) Acrolein	7.16	56	34469	114.48914 ppb		96
13) Acetone	7.27	43	22365	12.99757 ppb	#	84
14) Freon-113	7.46	101	138327	9.44782 ppb		97
15) 1,1-DCE	7.67	96	151407	8.83040 ppb		96
16) t-Butanol	7.76	59	10529	129.32077 ppb		93
17) Methyl Acetate	8.18	43	48755	9.36519 ppb		96
18) Iodomethane	8.16	142	102169	10.69989 ppb		90
19) Acrylonitrile	8.56	53	17916	9.49044 ppb		79
20) Methylene chloride	8.47	84	147951	9.04673 ppb		99
21) Carbon disulfide	8.56	76	154944	9.30051 ppb		100
22) Methyl t-butyl ether (MtBE	8.89	73	251165	9.69151 ppb		96
23) Trans-1,2-DCE	9.10	96	180083	9.05360 ppb		88
24) Diisopropyl Ether	9.75	45	553904	9.66784 ppb		94
25) 1,1-DCA	9.79	63	339012	9.97748 ppb		99
26) Vinyl Acetate	9.42	43	104836	9.85383 ppb	#	83
27) Ethyl tert Butyl Ether	10.45	59	395408	10.10888 ppb		99
28) MEK (2-Butanone)	10.44	43	71405	10.48433 ppb		99
29) Cis-1,2-DCE	10.82	96	187663	9.18084 ppb		97
30) 2,2-Dichloropropane	10.82	77	208247	8.55600 ppb		97
31) Chloroform	11.10	83	320091	9.79258 ppb		99
32) Bromochloromethane	11.32	128	58472	10.27501 ppb		98
34) 1,1,1-TCA	11.84	97	283983	9.55329 ppb		96
35) Cyclohexane	12.00	56	268948	9.71733 ppb		94
36) 1,1-Dichloropropene	12.10	75	240188	9.42981 ppb		99
37) 2,2,4-Trimethylpentane	12.18	57	414455	9.47423 ppb		98
39) Carbon Tetrachloride	12.30	117	199898	9.76577 ppb		98
40) Tert Amyl Methyl Ether	12.34	73	292021	9.98459 ppb		99
41) 1,2-DCA	12.38	62	161160	9.60516 ppb		100
42) Benzene	12.50	78	693647	9.47863 ppb		96
43) TCE	13.53	95	195399	9.64105 ppb		91

Algorithm Check: $\frac{(222193)(25)}{(600576)(0.92079)} (1) = 10.04166034$ ✓ ARS 12/7/11

(#) = qualifier out of range (m) = manual integration

Quantitation Report (Not Reviewed)

Data File : M:\CHICO\DATA\C111030\1030C28W.D Vial: 1
Acq On : 31 Oct 11 8:48 Operator: STC
Sample : 111030A LCS-1WC (SS) Inst : Chico
Misc : Water 10mLw/ IS&S:10-30/10-26-11 Multiplr: 1.00

Quant Time: Dec 2 11:35 2011 Quant Results File: CALLW.RES

Quant Method : M:\CHICO\DATA\C111030\CALLW.M (RTE Integrator)
Title : METHOD 8260
Last Update : Fri Dec 02 11:32:50 2011
Response via : Initial Calibration
DataAcq Meth : V8260

	Compound	R.T.	QIon	Response	Conc	Unit		Qvalue
44)	2-Pentanone	13.20	43	531259	125.31057	ppb		97
45)	1,2-Dichloropropane	13.76	63	167281	10.05687	ppb	#	94
46)	Bromodichloromethane	14.11	83	190021	9.99949	ppb	#	91
47)	Methyl Cyclohexane	13.82	83	224538	9.48040	ppb		99
48)	Dibromomethane	14.16	93	67527	10.15102	ppb		95
49)	2-Chloroethyl vinyl ether	14.57	63	42382	10.02259	ppb		95
50)	1-Bromo-2-chloroethane	14.88	63	147231	10.37368	ppb	#	79
51)	Cis-1,3-Dichloropropene	15.00	75	181879	10.03716	ppb		100
52)	Toluene	15.63	91	678338	9.39804	ppb		99
53)	Trans-1,3-Dichloropropene	15.80	75	126600	9.70449	ppb		98
54)	1,1,2-TCA	16.08	83	69681	9.91034	ppb		93
57)	1,2-EDB	17.33	107	79107	10.66417	ppb		94
58)	Tetrachloroethene	16.78	164	202010	10.08114	ppb		94
59)	1-Chlorohexane	17.70	91	232734	10.08580	ppb		97
60)	1,1,1,2-Tetrachloroethane	18.16	131	138607	11.04877	ppb		99
61)	m&p-Xylene	18.35	106	589147	19.90172	ppb		97
62)	o-Xylene	19.11	106	295217	10.36928	ppb		98
63)	Styrene	19.13	104	457607	10.64883	ppb		93
65)	2-Hexanone	16.11	43	38770	10.87089	ppb		95
66)	1,3-Dichloropropane	16.49	76	147530	10.08561	ppb		98
67)	Dibromochloromethane	16.97	129	105397	11.03714	ppb		82
68)	Chlorobenzene	18.10	112	437982	10.34243	ppb		97
69)	Ethylbenzene	18.22	91	794180	10.07104	ppb		100
70)	Bromoform	19.65	173	45131	9.31734	ppb		91
72)	MIBK (methyl isobutyl keto	14.68	43	59150	9.77763	ppb		87
73)	Isopropylbenzene	19.73	105	756513	9.80877	ppb		98
74)	1,1,2,2-Tetrachloroethane	19.90	83	66287	10.26718	ppb	#	74
75)	1,2,3-Trichloropropane	20.16	110	8565	11.81260	ppb		82
76)	t-1,4-Dichloro-2-Butene	20.23	53	14963	10.22116	ppb	#	92
77)	Bromobenzene	20.48	156	179052	10.06445	ppb		89
78)	n-Propylbenzene	20.44	91	900774	9.79012	ppb		100
79)	4-Ethyltoluene	20.63	105	593563	9.32229	ppb		97
80)	2-Chlorotoluene	20.74	91	598129	9.81561	ppb		98
81)	1,3,5-Trimethylbenzene	20.72	105	617840	9.86323	ppb		99
82)	4-Chlorotoluene	20.82	91	502123	9.56935	ppb		98
83)	Tert-Butylbenzene	21.36	119	672218	9.91209	ppb		97
84)	1,2,4-Trimethylbenzene	21.42	105	604092	9.23368	ppb		96
85)	Sec-Butylbenzene	21.76	105	823845	10.12964	ppb		96
86)	p-Isopropyltoluene	21.99	119	683604	9.81315	ppb		98
87)	Benzyl Chloride	22.42	91	81362	8.79846	ppb		94
88)	1,3-DCB	22.12	146	351790	9.66926	ppb		95
89)	1,4-DCB	22.30	146	335795	9.94420	ppb		96
90)	Hexachloroethane	23.59	117	92345	8.87470	ppb		87
91)	n-Butylbenzene	22.69	91	582962	9.59448	ppb		98
92)	1,2-DCB	22.93	146	292666	10.11316	ppb		97
93)	1,2-Dibromo-3-chloropropan	24.14	155	10559	10.20135	ppb		92
94)	1,2,4-Trichlorobenzene	25.59	180	213173	10.16551	ppb		98
95)	Hexachlorobutadiene	25.84	223	40176	10.54513	ppb		97
96)	Naphthalene	25.94	128	272964	10.54986	ppb		99
97)	1,2,3-Trichlorobenzene	26.29	180	172357	10.86589	ppb		99

(#) = qualifier out of range (m) = manual integration

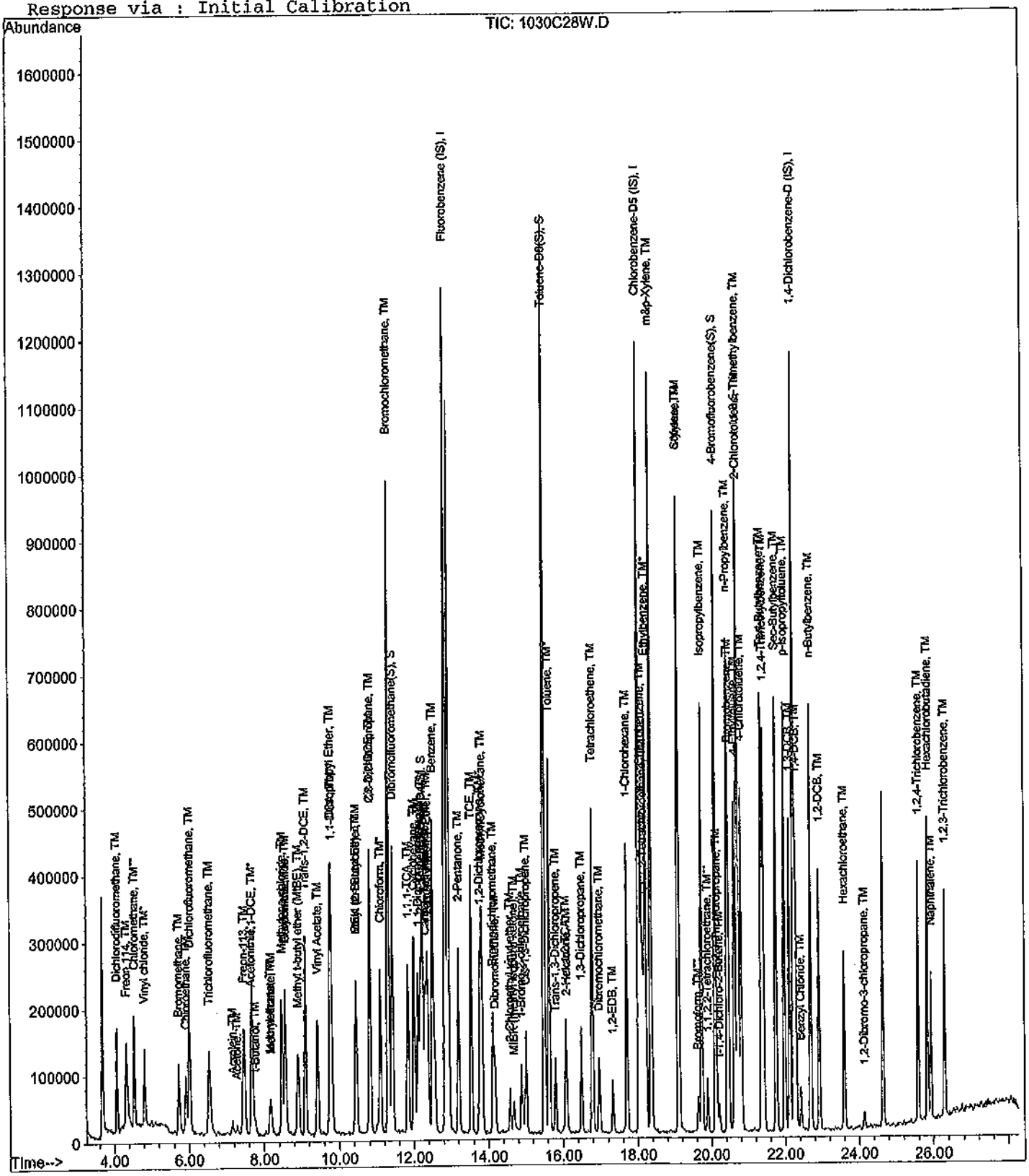
Quantitation Report

Data File : M:\CHICO\DATA\C111030\1030C28W.D
Acq On : 31 Oct 11 8:48
Sample : 111030A LCS-1WC (SS)
Misc : Water 10mLw/ IS&S:10-30/10-26-11

Vial: 1
Operator: STC
Inst : Chico
Multiplr: 1.00

Quant Time: Dec 2 11:35 2011 Quant Results File: CALLW.RES

Method : M:\CHICO\DATA\C111030\CALLW.M (RTE Integrator)
Title : METHOD 8260
Last Update : Fri Dec 02 11:32:50 2011
Response via : Initial Calibration



VOLATILE ORGANIC ANALYSIS BY
EPA METHOD 8260B

Form 7

Continuing Calibration

Lab Name: APPL, Inc.
Case No:
Matrix:

SDG No:
Date Analyzed: 10/31/2011
Instrument: Chico
Initial Cal. Date: 10/30/2011
Data File: 1030C27W.D

		Compound	MEAN	CCRF	%D		%Drift	
1	I	Fluorobenzene (IS)	ISTD			I		
2	TM	Dichlorodifluoromethane	0.9211	0.9607	4.3	TM		
3	TM	Freon 114	0.5769	0.6267	8.6	TM		
4	TM**	Chloromethane	1.141	1.117	2.2	TM**		
5	TM*	Vinyl chloride	0.7635	0.8485	11	TM*		
6	TML	1,3-Butadiene	0.0000	0.0009	0.00	TML		
7	TM	Bromomethane	0.5542	0.5584	0.75	TM		
8	TM	Chloroethane	0.6306	0.6136	2.7	TM		
9	TM	Dichlorofluoromethane	1.744	1.741	0.17	TM		
10	TM	Trichlorofluoromethane	1.035	1.087	5.1	TM		
11		Acetonitrile	0.0274	0.0255	6.7			
12	TM	Acrolein	0.0125	0.0117	6.4	TM		
13	TML	Acetone	0.1859	0.0914	51	TML	28	*
14	TML	Freon-113	0.5715	0.6270	9.7	TML	3.9	
15	TM*	1,1-DCE	0.7137	0.6880	3.6	TM*		
16	TM	t-Butanol	0.0034	0.0034	1.1	TM		
17	TML	Methyl Acetate	0.2927	0.2052	30	TML	5.2	
18	TML	Iodomethane	0.3500	0.4294	23	TML	7.7	
19	TML	Acrylonitrile	0.0764	0.0791	3.5	TML	0.85	
20	TM	Methylene chloride	0.6808	0.6482	4.8	TM		
21	TM	Carbon disulfide	0.6935	0.6961	0.37	TM		
22	TM	Methyl t-butyl ether (MtBE)	1.079	1.106	2.6	TM		
23	TM	Trans-1,2-DCE	0.8280	0.7771	6.1	TM		
24	TM	Diisopropyl Ether	2.385	2.416	1.3	TM		
25	TM**	1,1-DCA	1.414	1.481	4.7	TM**		
26	TML	Vinyl Acetate	0.5623	0.4689	17	TML	7.0	
27	TM	Ethyl tert Butyl Ether	1.628	1.743	7.1	TM		
28	TML	MEK (2-Butanone)	0.3591	0.2865	20	TML	0.79	
29	TM	Cis-1,2-DCE	0.8509	0.8285	2.6	TM		
30	TM	2,2-Dichloropropane	1.013	0.9291	8.3	TM		
31	TM*	Chloroform	1.361	1.374	1.0	TM*		
32	TM	Bromochloromethane	0.2369	0.2573	8.6	TM		
33	S	Dibromofluoromethane(S)	0.6660	0.6943	4.3	S		
34	TM	1,1,1-TCA	1.237	1.277	3.2	TM		
35	TM	Cyclohexane	1.152	1.162	0.83	TM		
36	TM	1,1-Dichloropropene	1.060	1.091	2.9	TM		
37	TML	2,2,4-Trimethylpentane	2.316	1.777	23	TML	2.1	
38	S	1,2-DCA-D4(S)	0.5928	0.5850	1.3	S		
39	TM	Carbon Tetrachloride	0.8521	0.9156	7.5	TM		
40	TM	Tert Amyl Methyl Ether	1.217	1.268	4.1	TM		
		Average			7.7			

VOLATILE ORGANIC ANALYSIS BY
EPA METHOD 8260B

Form 7

Continuing Calibration

Lab Name: APPL, Inc.
Case No:
Matrix: 0

SDG No:
Date Analyzed: 10/31/2011
Instrument: Chico
Cal. Date: 10/30/2011
Data File: 1030C27W.D

		Compound	MEAN	CCRF	%D		%Drift
41	TM	1,2-DCA	0.6984	0.7162	2.5	TM	
42	TM	Benzene	3.046	3.004	1.4	TM	
43	TM	TCE	0.8437	0.8976	6.4	TM	
44	TM	2-Pentanone	0.1765	0.1801	2.0	TM	
45	TM*	1,2-Dichloropropane	0.6924	0.7108	2.7	TM*	
46	TM	Bromodichloromethane	0.7910	0.8325	5.2	TM	
47	TM	Methyl Cyclohexane	0.9859	0.9762	0.98	TM	
48	TM	Dibromomethane	0.2769	0.3053	10	TM	
49	TM	2-Chloroethyl vinyl ether	0.1760	0.1791	1.8	TM	
50	TM	1-Bromo-2-chloroethane	0.5908	0.6158	4.2	TM	
51	TM	Cis-1,3-Dichloropropene	0.7543	0.7871	4.4	TM	
52	TM*	Toluene	3.005	3.021	0.56	TM*	
53	TM	Trans-1,3-Dichloropropene	0.5430	0.5626	3.6	TM	
54	TM	1,1,2-TCA	0.2927	0.3204	9.5	TM	
55	I	Chlorobenzene-D5 (IS)	ISTD			I	
56	S	Toluene-D8(S)	3.518	3.735	6.2	S	
57	TM	1,2-EDB	0.4758	0.5192	9.1	TM	
58	TM	Tetrachloroethene	1.285	1.322	2.9	TM	
59	TM	1-Chlorohexane	1.480	1.498	1.2	TM	
60	TM	1,1,1,2-Tetrachloroethane	0.8047	0.9010	12	TM	
61	TM	m&p-Xylene	1.899	1.872	1.4	TM	
62	TM	o-Xylene	1.826	1.841	0.81	TM	
63	TM	Styrene	2.756	2.893	5.0	TM	
64	S	4-Bromofluorobenzene(S)	1.260	1.340	6.3	S	
65	TM	2-Hexanone	0.2288	0.2478	8.3	TM	
66	TM	1,3-Dichloropropane	0.9383	0.9771	4.1	TM	
67	TM	Dibromochloromethane	0.6125	0.6543	6.8	TM	
68	TM**	Chlorobenzene	2.716	2.717	0.01	TM**	
69	TM*	Ethylbenzene	5.058	5.085	0.53	TM*	
70	TM**L	Bromoform	0.2607	0.2986	15	TM**L	4.2
71	I	1,4-Dichlorobenzene-D (IS)	ISTD			I	
72	TM	MIBK (methyl isobutyl ketone)	0.7107	0.6949	2.2	TM	
73	TM	Isopropylbenzene	9.061	9.158	1.1	TM	
74	TM**	1,1,2,2-Tetrachloroethane	0.7585	0.7768	2.4	TM**	
75	TML	1,2,3-Trichloropropane	0.0967	0.1005	4.0	TML	18
76	TM	t-1,4-Dichloro-2-Butene	0.1720	0.1640	4.7	TM	
77	TM	Bromobenzene	2.090	2.096	0.30	TM	
78	TM	n-Propylbenzene	10.8	11.1	2.9	TM	
79	TM	4-Ethyltoluene	7.480	7.378	1.4	TM	
80	TM	2-Chlorotoluene	7.159	7.296	1.9	TM	
		Average			4.1		

VOLATILE ORGANIC ANALYSIS BY
EPA METHOD 8260B

Form 7

Continuing Calibration

Lab Name: APPL, Inc.
Case No:
Matrix: 0

SDG No:
Date Analyzed: 10/31/2011
Instrument: Chico
Cal. Date: 10/30/2011
Data File: 1030C27W.D

		Compound	MEAN	CCRF	%D		%Drift
81	TM	1,3,5-Trimethylbenzene	7.359	7.417	0.79	TM	
82	TM	4-Chlorotoluene	6.164	6.076	1.4	TM	
83	TM	Tert-Butylbenzene	7.967	8.229	3.3	TM	
84	TM	1,2,4-Trimethylbenzene	7.686	7.406	3.6	TM	
85	TM	Sec-Butylbenzene	9.555	9.767	2.2	TM	
86	TM	p-Isopropyltoluene	8.184	8.300	1.4	TM	
87	TM	Benzyl Chloride	1.086	0.9882	9.0	TM	
88	TM	1,3-DCB	4.274	4.305	0.72	TM	
89	TM	1,4-DCB	3.967	4.065	2.5	TM	
90	TML	Hexachloroethane	1.021	1.118	9.5	TML	9.2
91	TM	n-Butylbenzene	7.138	6.965	2.4	TM	
92	TM	1,2-DCB	3.400	3.527	3.7	TM	
93	TML	1,2-Dibromo-3-chloropropane	0.1148	0.1161	1.2	TML	3.7
94	TM	1,2,4-Trichlorobenzene	2.464	2.540	3.1	TM	
95	TM	Hexachlorobutadiene	0.4476	0.4532	1.2	TM	
96	TM	Naphthalene	3.040	3.215	5.8	TM	
97	TM	1,2,3-Trichlorobenzene	1.864	1.911	2.5	TM	
98							
99							
100							
101							
102							
103							
104							
105							
106							
107							
108							
109							
110							
111							
112							
113							
114							
115							
116							
117							
118							
119							
120							
		Average			3.2		

Quantitation Report (Not Reviewed)

Data File : M:\CHICO\DATA\C111030\1030C27W.D Vial: 1
Acq On : 31 Oct 11 8:05 Operator: STC
Sample : Voc Std 10-30-11@10ug/L Inst : Chico
Misc : Water 10mLw/ IS&S:10-30/10-26-11 Multiplr: 1.00

Quant Time: Dec 2 11:35 2011 Quant Results File: CALLW.RES

Quant Method : M:\CHICO\DATA\C111030\CALLW.M (RTE Integrator)
Title : METHOD 8260
Last Update : Fri Dec 02 11:32:50 2011
Response via : Initial Calibration
DataAcq Meth : V8260

Internal Standards	R.T.	QIon	Response	Conc	Units		Dev(Min)
1) Fluorobenzene (IS)	12.84	96	583168	25.00000	ppb		0.00
55) Chlorobenzene-D5 (IS)	18.03	117	397504	25.00000	ppb		0.00
71) 1,4-Dichlorobenzene-D (IS)	22.24	152	209984	25.00000	ppb		0.00
System Monitoring Compounds							
33) Dibromofluoromethane(S)	11.43	111	404916	26.06515	ppb		0.00
Spiked Amount 25.097			Recovery = 103.856%				
38) 1,2-DCA-D4(S)	12.23	65	341171	24.67132	ppb		0.00
Spiked Amount 24.225			Recovery = 101.840%				
56) Toluene-D8(S)	15.50	98	1484769	26.54660	ppb		0.00
Spiked Amount 25.808			Recovery = 102.862%				
64) 4-Bromofluorobenzene(S)	20.11	95	532764	26.58449	ppb		0.00
Spiked Amount 25.459			Recovery = 104.417%				
Target Compounds							Qvalue
2) Dichlorodifluoromethane	4.06	85	224089	10.42966	ppb		97
3) Freon 114	4.33	85	146183	10.86336	ppb		98
4) Chloromethane	4.55	50	260469	9.78278	ppb		98
5) Vinyl chloride	4.82	62	197919	11.11320	ppb		95
7) Bromomethane	5.72	94	130249	10.07502	ppb		89
8) Chloroethane	5.91	64	143140	9.73031	ppb		99
9) Dichlorofluoromethane	6.00	67	406047	9.98304	ppb		100
10) Trichlorofluoromethane	6.52	101	253658	10.50908	ppb		97
11) Acetonitrile	7.65	41	74497	116.61139	ug/l		100
12) Acrolein	7.15	56	34193	116.96262	ppb		94
13) Acetone	7.29	43	21316	12.75773	ppb	##	77
14) Freon-113	7.46	101	146261	10.39098	ppb		94
15) 1,1-DCE	7.68	96	160490	9.63955	ppb		97
16) t-Butanol	7.76	59	9773	123.61847	ppb		100
17) Methyl Acetate	8.19	43	47860	9.47524	ppb		98
18) Iodomethane	8.17	142	100159	10.76585	ppb	#	89
19) Acrylonitrile	8.55	53	18444	10.08467	ppb		84
20) Methylene chloride	8.48	84	151205	9.52169	ppb		86
21) Carbon disulfide	8.55	76	162368	10.03707	ppb		98
22) Methyl t-butyl ether (MtBE	8.90	73	258098	10.25631	ppb		95
23) Trans-1,2-DCE	9.09	96	181275	9.38557	ppb		90
24) Diisopropyl Ether	9.75	45	563558	10.12996	ppb		96
25) 1,1-DCA	9.79	63	345459	10.47072	ppb		97
26) Vinyl Acetate	9.42	43	109376	10.70268	ppb	#	83
27) Ethyl tert Butyl Ether	10.45	59	406696	10.70784	ppb		94
28) MEK (2-Butanone)	10.43	43	66821	10.07941	ppb		96
29) Cis-1,2-DCE	10.81	96	193255	9.73664	ppb		97
30) 2,2-Dichloropropane	10.81	77	216729	9.17030	ppb		95
31) Chloroform	11.09	83	320623	10.10165	ppb		99
32) Bromochloromethane	11.31	128	60012	10.86042	ppb		81
34) 1,1,1-TCA	11.83	97	297856	10.31908	ppb		97
35) Cyclohexane	12.01	56	270975	10.08282	ppb		92
36) 1,1-Dichloropropene	12.10	75	254512	10.29044	ppb		100
37) 2,2,4-Trimethylpentane	12.18	57	414560	9.79424	ppb		98
39) Carbon Tetrachloride	12.30	117	213577	10.74550	ppb		94
40) Tert Amyl Methyl Ether	12.34	73	295693	10.41194	ppb		98

(#) = qualifier out of range (m) = manual integration

Quantitation Report (Not Reviewed)

Data File : M:\CHICO\DATA\C111030\1030C27W.D Vial: 1
Acq On : 31 Oct 11 8:05 Operator: STC
Sample : Voc Std 10-30-11@10ug/L Inst : Chico
Misc : Water 10mLw/ IS&S:10-30/10-26-11 Multiplr: 1.00

Quant Time: Dec 2 11:35 2011 Quant Results File: CALLW.RES

Quant Method : M:\CHICO\DATA\C111030\CALLW.M (RTE Integrator)
Title : METHOD 8260
Last Update : Fri Dec 02 11:32:50 2011
Response via : Initial Calibration
DataAcq Meth : V8260

	Compound	R.T.	QIon	Response	Conc Unit		Qvalue
41)	1,2-DCA	12.37	62	167070	10.25463 ppb		96
42)	Benzene	12.50	78	700767	9.86177 ppb		97
43)	TCE	13.53	95	209384	10.63947 ppb		91
44)	2-Pentanone	13.20	43	525053	127.54364 ppb		98
45)	1,2-Dichloropropane	13.76	63	165813	10.26618 ppb	#	93
46)	Bromodichloromethane	14.12	83	194189	10.52386 ppb	#	92
47)	Methyl Cyclohexane	13.82	83	227722	9.90185 ppb		99
48)	Dibromomethane	14.17	93	71228	11.02699 ppb		99
49)	2-Chloroethyl vinyl ether	14.57	63	41783	10.17589 ppb		97
50)	1-Bromo-2-chloroethane	14.88	63	143648	10.42335 ppb		87
51)	Cis-1,3-Dichloropropene	15.01	75	183608	10.43504 ppb		97
52)	Toluene	15.63	91	704784	10.05591 ppb		96
53)	Trans-1,3-Dichloropropene	15.80	75	131238	10.36031 ppb		91
54)	1,1,2-TCA	16.09	83	74743	10.94760 ppb		93
57)	1,2-EDB	17.33	107	82556	10.91231 ppb		91
58)	Tetrachloroethene	16.79	164	210273	10.28907 ppb		91
59)	1-Chlorohexane	17.70	91	238146	10.11928 ppb		96
60)	1,1,1,2-Tetrachloroethane	18.16	131	143265	11.19759 ppb		93
61)	m&p-Xylene	18.35	106	595342	19.71920 ppb		97
62)	o-Xylene	19.11	106	292710	10.08093 ppb		98
63)	Styrene	19.12	104	460015	10.49632 ppb		97
65)	2-Hexanone	16.10	43	39396	10.83121 ppb	#	75
66)	1,3-Dichloropropane	16.49	76	155367	10.41445 ppb		95
67)	Dibromochloromethane	16.97	129	104038	10.68258 ppb		89
68)	Chlorobenzene	18.10	112	431948	10.00123 ppb		95
69)	Ethylbenzene	18.22	91	808497	10.05286 ppb		95
70)	Bromoform	19.64	173	47481	9.57573 ppb	#	74
72)	MIBK (methyl isobutyl keto	14.68	43	58368	9.77776 ppb		91
73)	Isopropylbenzene	19.74	105	769191	10.10690 ppb		99
74)	1,1,2,2-Tetrachloroethane	19.90	83	65243	10.24099 ppb		86
75)	1,2,3-Trichloropropane	20.15	110	8442	11.79921 ppb		79
76)	t-1,4-Dichloro-2-Butene	20.23	53	13773	9.53445 ppb	#	90
77)	Bromobenzene	20.48	156	176075	10.02984 ppb		93
78)	n-Propylbenzene	20.44	91	934437	10.29219 ppb		100
79)	4-Ethyltoluene	20.64	105	619715	9.86355 ppb		96
80)	2-Chlorotoluene	20.73	91	612789	10.19105 ppb		97
81)	1,3,5-Trimethylbenzene	20.71	105	623010	10.07914 ppb		97
82)	4-Chlorotoluene	20.82	91	510371	9.85698 ppb		98
83)	Tert-Butylbenzene	21.36	119	691172	10.32825 ppb		98
84)	1,2,4-Trimethylbenzene	21.42	105	622056	9.63577 ppb		97
85)	Sec-Butylbenzene	21.76	105	820355	10.22199 ppb		99
86)	p-Isopropyltoluene	21.99	119	697110	10.14123 ppb		99
87)	Benzyl Chloride	22.42	91	83004	9.09640 ppb		99
88)	1,3-DCB	22.13	146	361602	10.07224 ppb		95
89)	1,4-DCB	22.30	146	341438	10.24691 ppb		96
90)	Hexachloroethane	23.60	117	93885	9.08419 ppb		96
91)	n-Butylbenzene	22.70	91	585015	9.75739 ppb		98
92)	1,2-DCB	22.93	146	296267	10.37488 ppb		95
93)	1,2-Dibromo-3-chloropropan	24.14	155	9755	9.63316 ppb		88
94)	1,2,4-Trichlorobenzene	25.59	180	213308	10.30836 ppb		97
95)	Hexachlorobutadiene	25.84	223	38064	10.12477 ppb		98

(#) = qualifier out of range (m) = manual integration

Quantitation Report (Not Reviewed)

Data File : M:\CHICO\DATA\C111030\1030C27W.D Vial: 1
Acq On : 31 Oct 11 8:05 Operator: STC
Sample : Voc Std 10-30-11@10ug/L Inst : Chico
Misc : Water 10mLw/ IS&S:10-30/10-26-11 Multiplr: 1.00

Quant Time: Dec 2 11:35 2011 Quant Results File: CALLW.RES

Quant Method : M:\CHICO\DATA\C111030\CALLW.M (RTE Integrator)
Title : METHOD 8260
Last Update : Fri Dec 02 11:32:50 2011
Response via : Initial Calibration
DataAcq Meth : V8260

Compound	R.T.	QIon	Response	Conc Unit	Qvalue
96) Naphthalene	25.94	128	270062	10.57768 ppb	99
97) 1,2,3-Trichlorobenzene	26.29	180	160478	10.25268 ppb	93

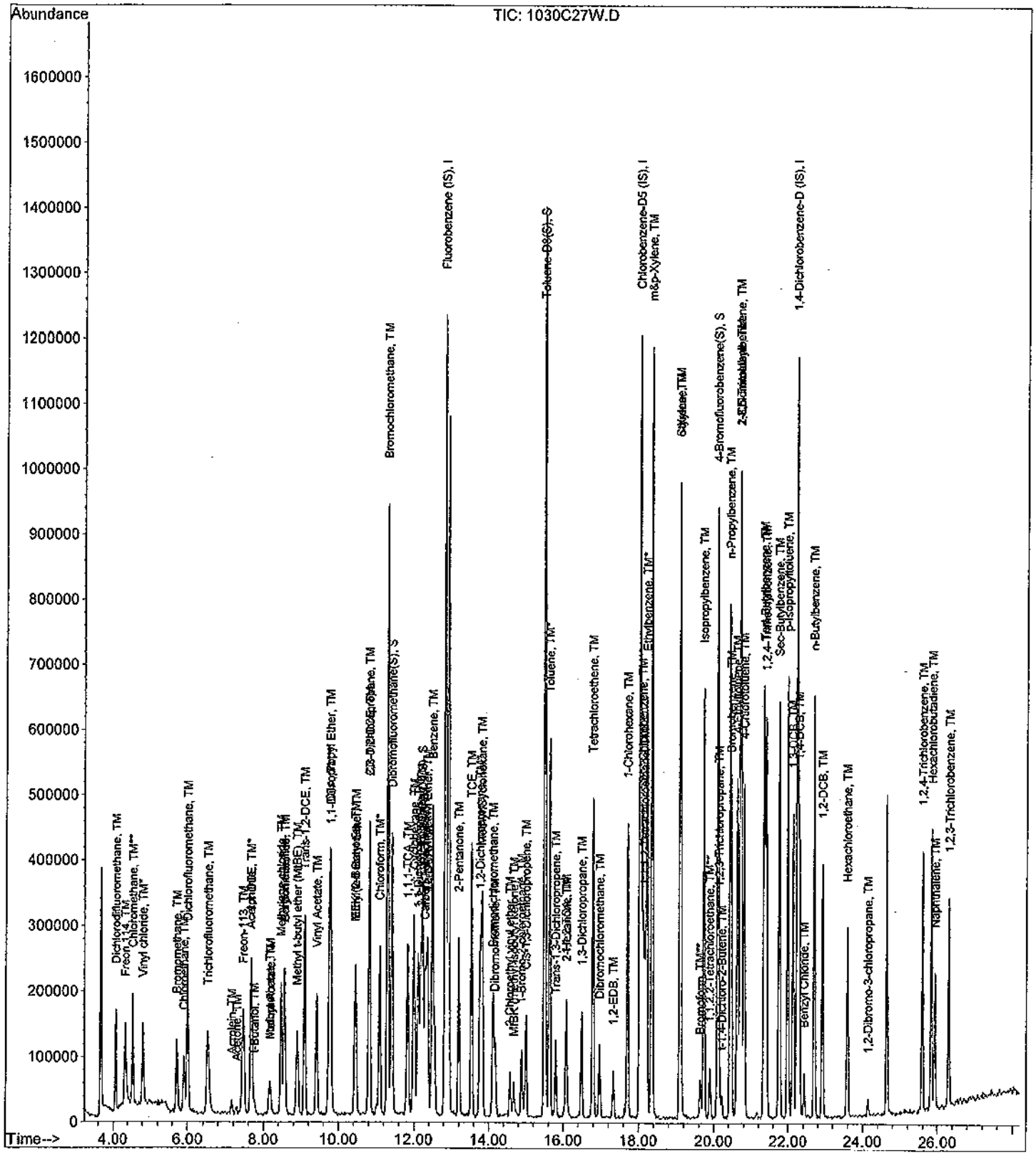
(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : M:\CHICO\DATA\C111030\1030C27W.D Vial: 1
Acq On : 31 Oct 11 8:05 Operator: STC
Sample : Voc Std 10-30-11@10ug/L Inst : Chico
Misc : Water 10mLw/ IS&S:10-30/10-26-11 Multiplr: 1.00

Quant Time: Dec 2 11:35 2011 Quant Results File: CALLW.RES

Method : M:\CHICO\DATA\C111030\CALLW.M (RTE Integrator)
Title : METHOD 8260
Last Update : Fri Dec 02 11:32:50 2011
Response via : Initial Calibration



VOLATILE ORGANIC ANALYSIS BY
EPA METHOD 8260B

Form 6
Initial Calibration

Lab Name: APPL, Inc. SDG No:
Case No: Initial Cal. Date: 10/30/2011
Matrix: Instrument: Chico

Initials:

		Compound	0.5	1	5	10	20	40	100				Avg	%RSD		
			1030C05W.D	1030C06W.D	1030C07W.D	1030C08W.D	1030C09W.D	1030C10W.D	1030C11W.D							
1	I	Fluorobenzene (IS)	ISTD													
2	TMHBL	Gasoline	17.8	8.144	5.154	3.050	2.484	2.343	2.274				5.9	96	TMHBL	1.000
3	I	Chlorobenzene-D5 (IS)	ISTD													
4	I	1,4-Dichlorobenzene-D (IS)	ISTD													
5																
6																
7																
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Quantitation Report (QT Reviewed)

Data File : M:\CHICO\DATA\C111030\1030C05W.D
Acq On : 30 Oct 11 16:17
Sample : Vol Std 10-30-11@20ug/L
Misc : Water 10mLw/ IS:10-30-11

Vial: 1
Operator: STC
Inst : Chico
Multiplr: 1.00

Quant Time: Nov 3 10:29 2011 Quant Results File: CGAS.RES

Quant Method : M:\CHICO\DATA\C111030\CGAS.M (RTE Integrator)
Title : METHOD 8260B
Last Update : Mon Oct 31 09:32:18 2011
Response via : Initial Calibration
DataAcq Meth : V8260

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Fluorobenzene (IS)	12.86	TIC	1064868	25.00000	ppb	0.02
3) Chlorobenzene-D5 (IS)	18.05	TIC	1075283	25.00000	ppb	0.01
4) 1,4-Dichlorobenzene-D (IS)	22.26	TIC	1031464	25.00000	ppb	0.02
System Monitoring Compounds						
Target Compounds						Qvalue
2) Gasoline	18.05	TIC	15186538m	62.79631	ppb	100

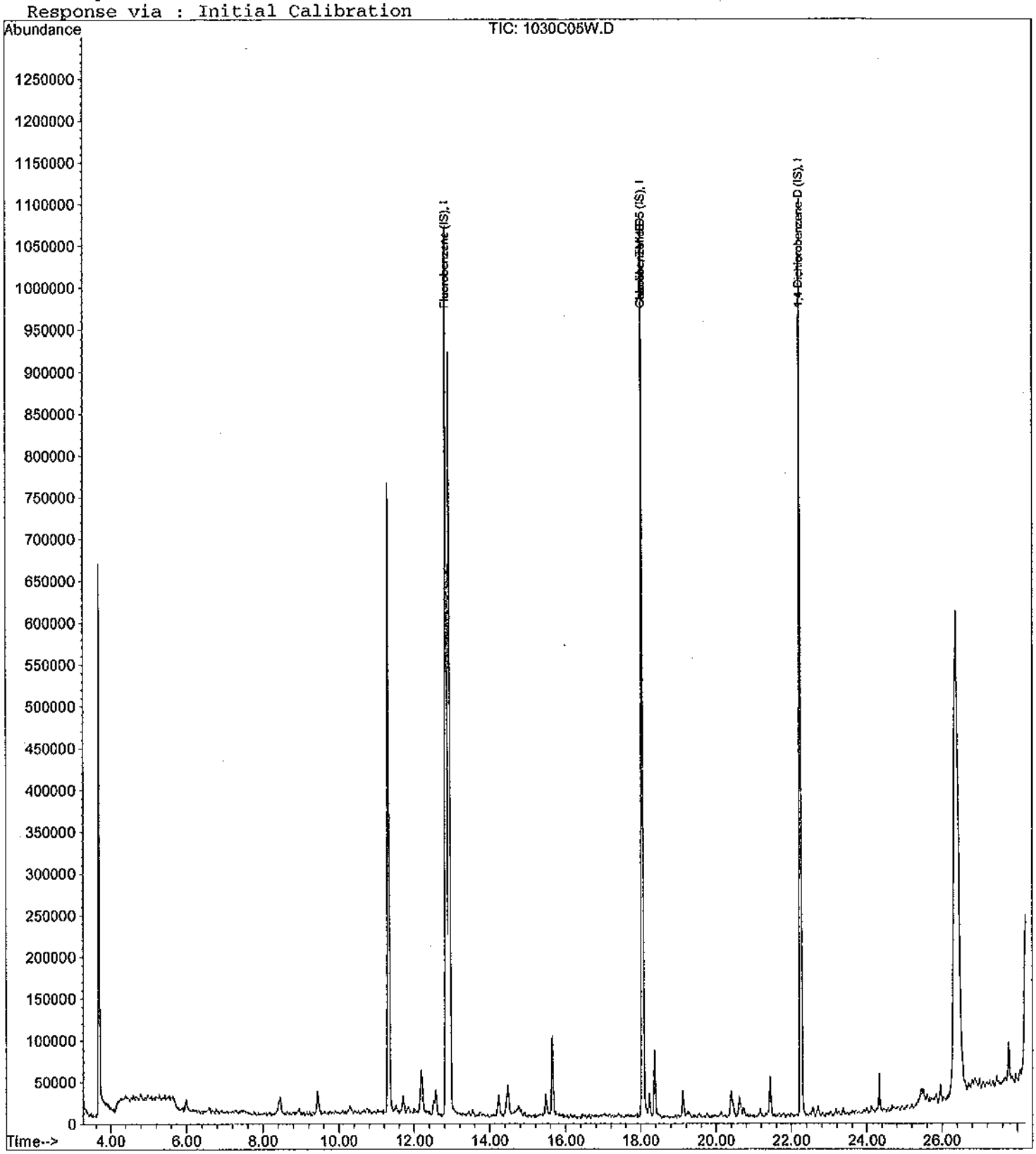
(#) = qualifier out of range (m) = manual integration
1030C05W.D CGAS.M Thu Nov 03 11:11:58 2011

Quantitation Report

Data File : M:\CHICO\DATA\C111030\1030C05W.D Vial: 1
Acq On : 30 Oct 11 16:17 Operator: STC
Sample : Vol Std 10-30-11@20ug/L Inst : Chico
Misc : Water 10mLw/ IS:10-30-11 Multiplr: 1.00

Quant Time: Nov 3 10:29 2011 Quant Results File: CGAS.RES

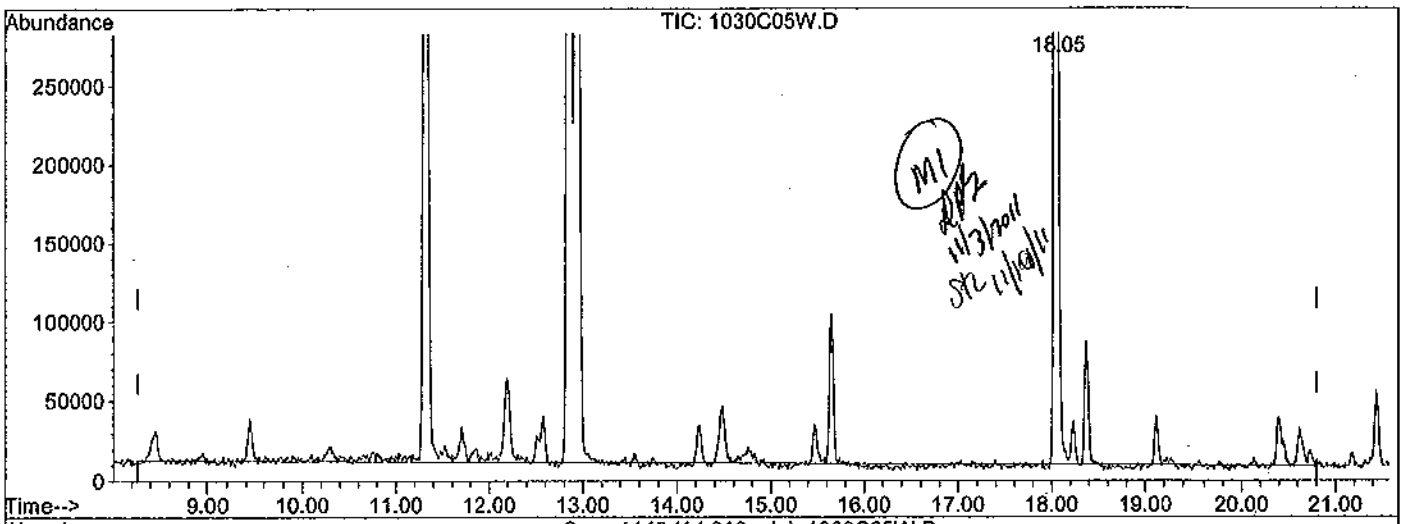
Method : M:\CHICO\DATA\C111030\CGAS.M (RTE Integrator)
Title : METHOD 8260B
Last Update : Thu Nov 03 10:47:02 2011
Response via : Initial Calibration

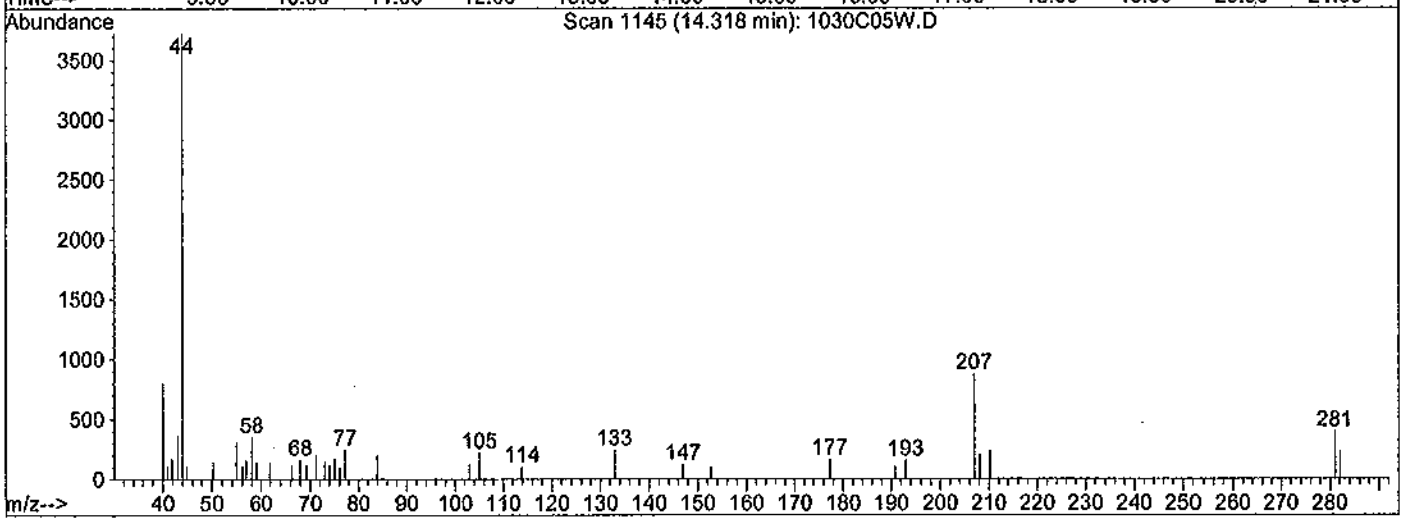


Quantitation Report

Data File : M:\CHICO\DATA\C111030\1030C05W.D Vial: 1
Acq On : 30 Oct 11 16:17 Operator: STC
Sample : Vol Std 10-30-11@20ug/L Inst : Chico
Misc : Water 10mLw/ IS:10-30-11 Multiplr: 1.00
Quant Time: Oct 31 9:32 2011 Quant Results File: temp.res

Method : M:\CHICO\DATA\C111030\CGAS.M (RTE Integrator)
Title : METHOD 8260B
Last Update : Mon Oct 31 09:35:31 2011
Response via : Multiple Level Calibration





TIC: 1030C05W.D

(2) Gasoline (TMHB)

14.31min 48.1330ppb m

response 11640400

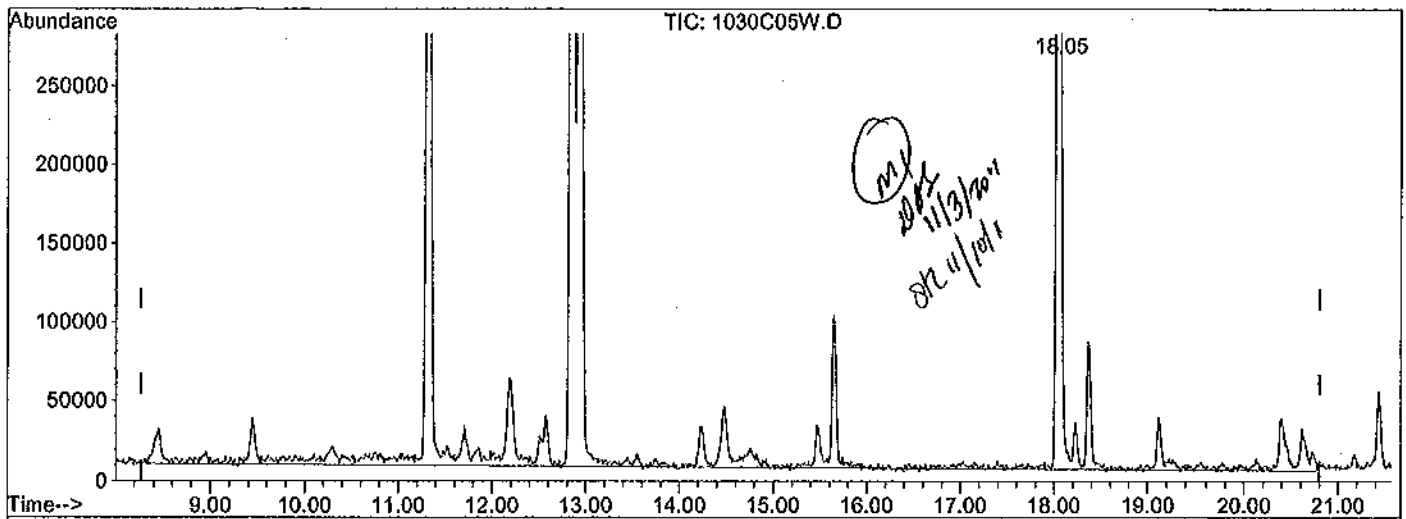
Ion	Exp%	Act%
TIC	100	100
0.00	0.00	0.82#
0.00	0.00	2.44#
0.00	0.00	0.00

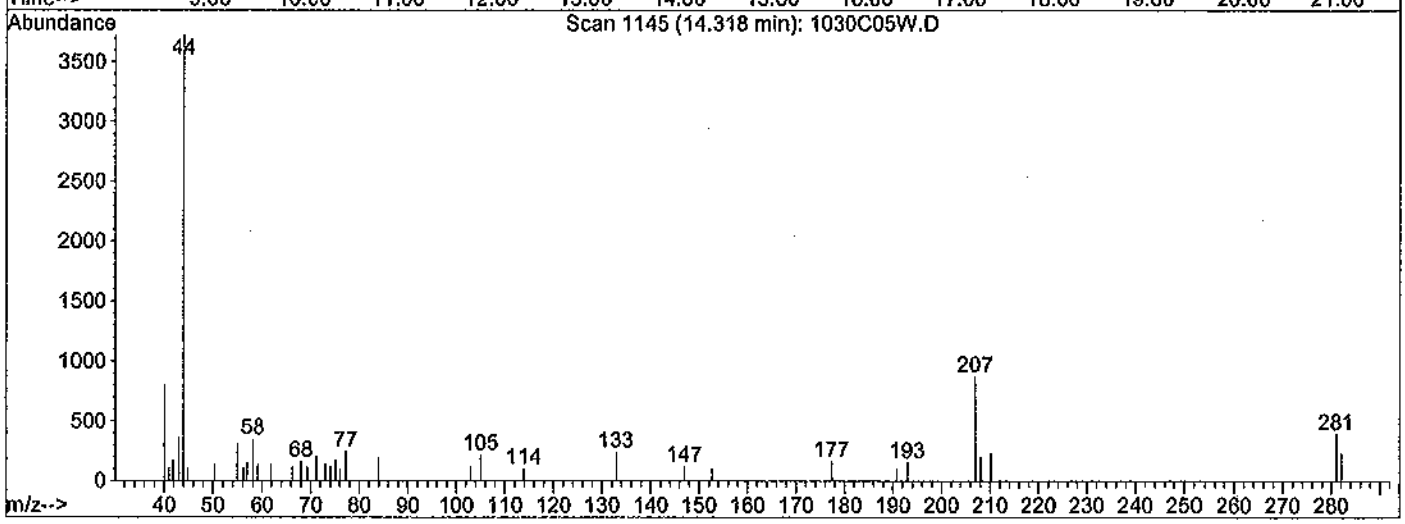
1030C05W.D CGAS.M Thu Nov 03 10:29:03 2011

Quantitation Report

Data File : M:\CHICO\DATA\C111030\1030C05W.D Vial: 1
Acq On : 30 Oct 11 16:17 Operator: STC
Sample : Vol Std 10-30-11@20ug/L Inst : Chico
Misc : Water 10mLw/ IS:10-30-11 Multiplr: 1.00
Quant Time: Nov 3 10:29 2011 Quant Results File: temp.res

Method : M:\CHICO\DATA\C111030\CGAS.M (RTE Integrator)
Title : METHOD 8260B
Last Update : Mon Oct 31 09:35:31 2011
Response via : Multiple Level Calibration





TIC: 1030C05W.D

(2) Gasoline (TMHB)

18.05min 62.7963ppb m

response 15186538

Ion	Exp%	Act%
TIC	100	100
0.00	0.00	0.63#
0.00	0.00	1.87#
0.00	0.00	0.00

Quantitation Report (QT Reviewed)

Data File : M:\CHICO\DATA\C111030\1030C06W.D Vial: 1
Acq On : 30 Oct 11 17:00 Operator: STC
Sample : Vol Std 10-30-11@50ug/L Inst : Chico
Misc : Water 10mLw/ IS:10-30-11 Multiplr: 1.00

Quant Time: Nov 3 10:30 2011 Quant Results File: CGAS.RES

Quant Method : M:\CHICO\DATA\C111030\CGAS.M (RTE Integrator)
Title : METHOD 8260B
Last Update : Mon Oct 31 09:32:18 2011
Response via : Initial Calibration
DataAcq Meth : V8260

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Fluorobenzene (IS)	12.85	TIC	1074535	25.00000	ppb	0.00
3) Chlorobenzene-D5 (IS)	18.04	TIC	1105653	25.00000	ppb	0.00
4) 1,4-Dichlorobenzene-D (IS)	22.24	TIC	1049854	25.00000	ppb	0.00

System Monitoring Compounds

Target Compounds						Qvalue
2) Gasoline	18.04	TIC	17501250m	71.71659	ppb	100

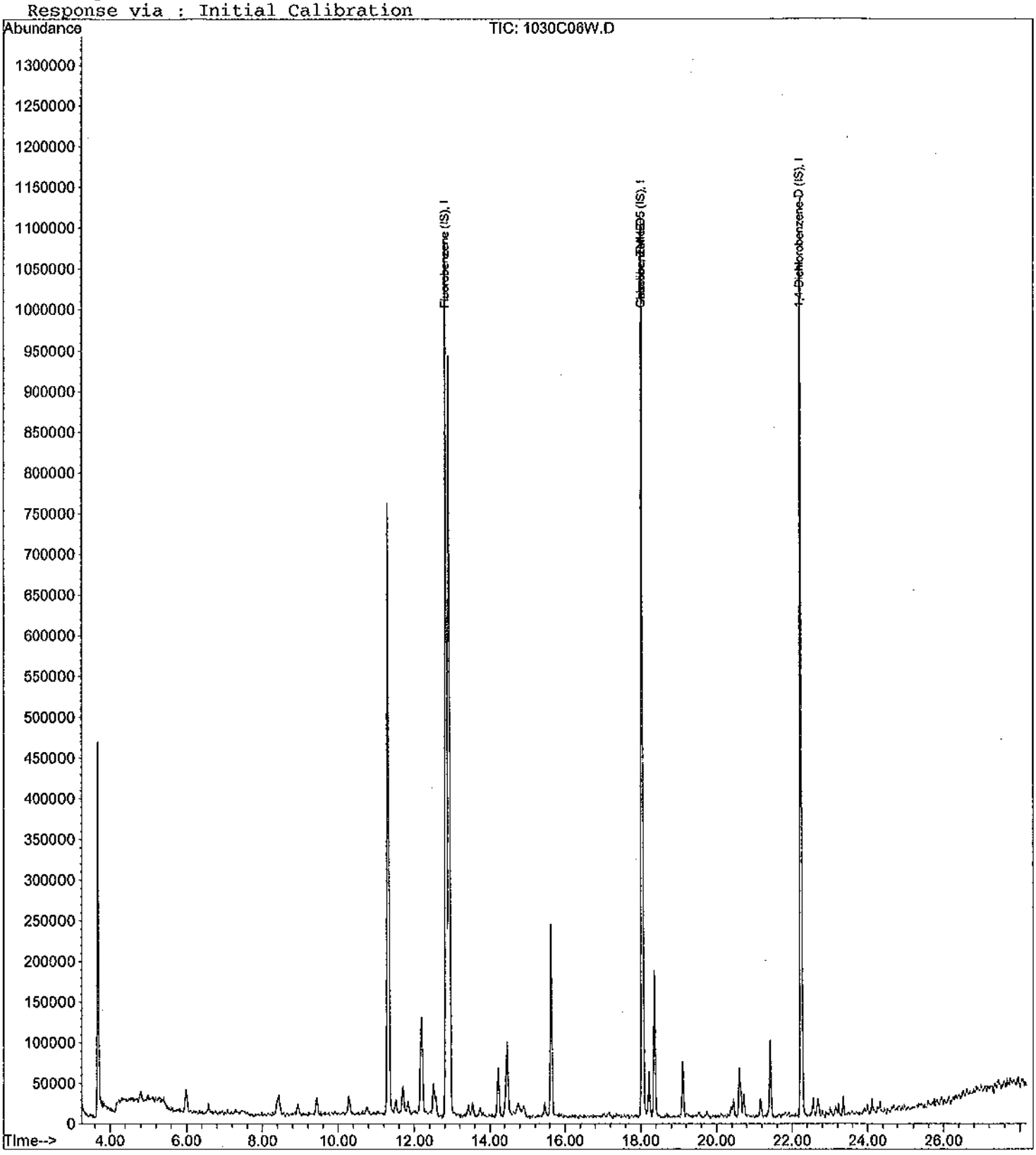
(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : M:\CHICO\DATA\C111030\1030C06W.D Vial: 1
Acq On : 30 Oct 11 17:00 Operator: STC
Sample : Vol Std 10-30-11@50ug/L Inst : Chico
Misc : Water 10mLw/ IS:10-30-11 Multiplr: 1.00

Quant Time: Nov 3 10:30 2011 Quant Results File: CGAS.RES

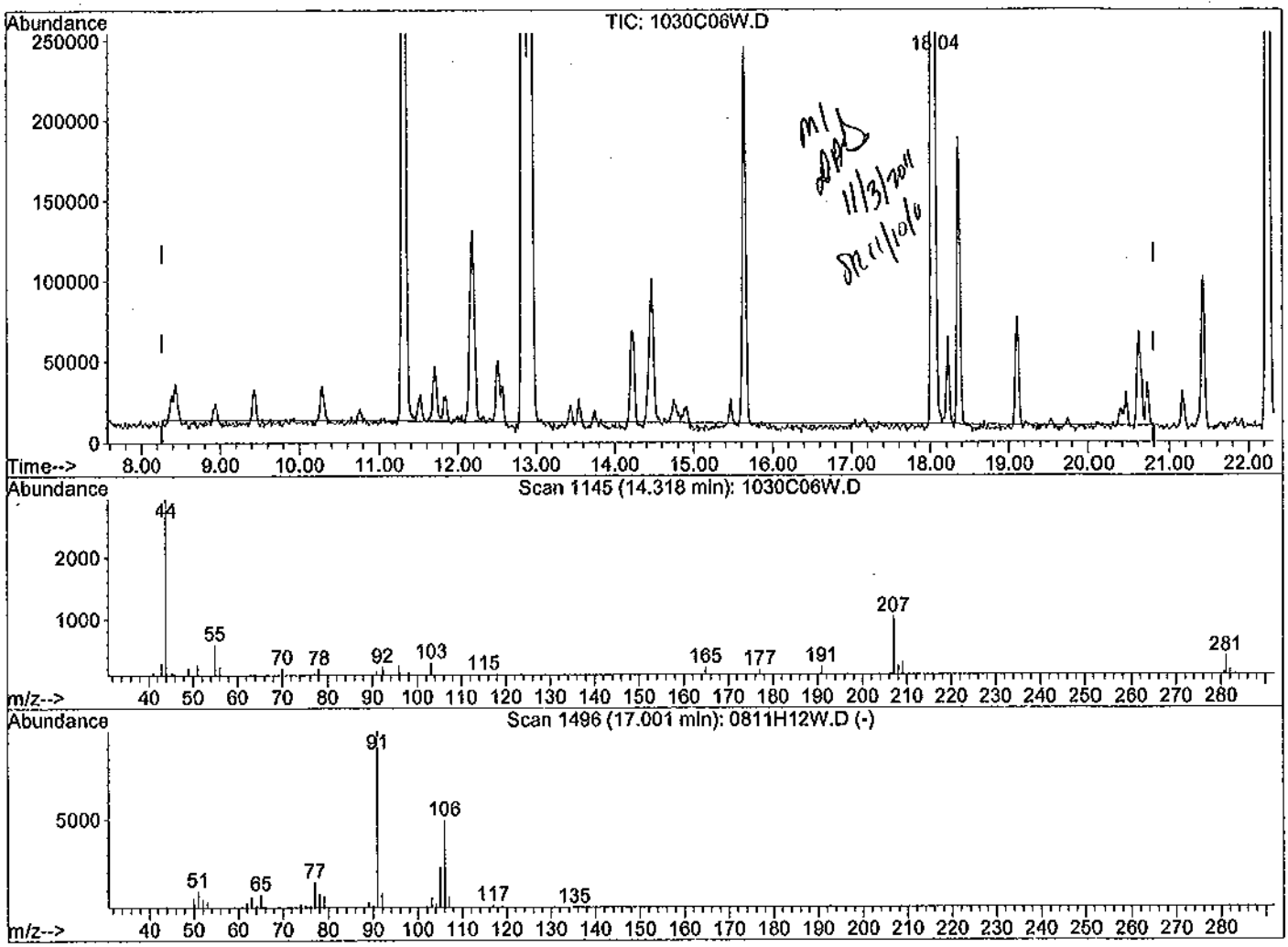
Method : M:\CHICO\DATA\C111030\CGAS.M (RTE Integrator)
Title : METHOD 8260B
Last Update : Thu Nov 03 10:47:02 2011
Response via : Initial Calibration



Quantitation Report

Data File : M:\CHICO\DATA\C111030\1030C06W.D Vial: 1
Acq On : 30 Oct 11 17:00 Operator: STC
Sample : Vol Std 10-30-11@50ug/L Inst : Chico
Misc : Water 10mLw/ IS:10-30-11 Multiplr: 1.00
Quant Time: Oct 31 9:32 2011 Quant Results File: temp.res

Method : M:\CHICO\DATA\C111030\CGAS.M (RTE Integrator)
Title : METHOD 8260B
Last Update : Mon Oct 31 09:35:31 2011
Response via : Multiple Level Calibration



TIC: 1030C06W.D

(2) Gasoline (TMHB)

14.31min 58.1426ppb m

response 14188741

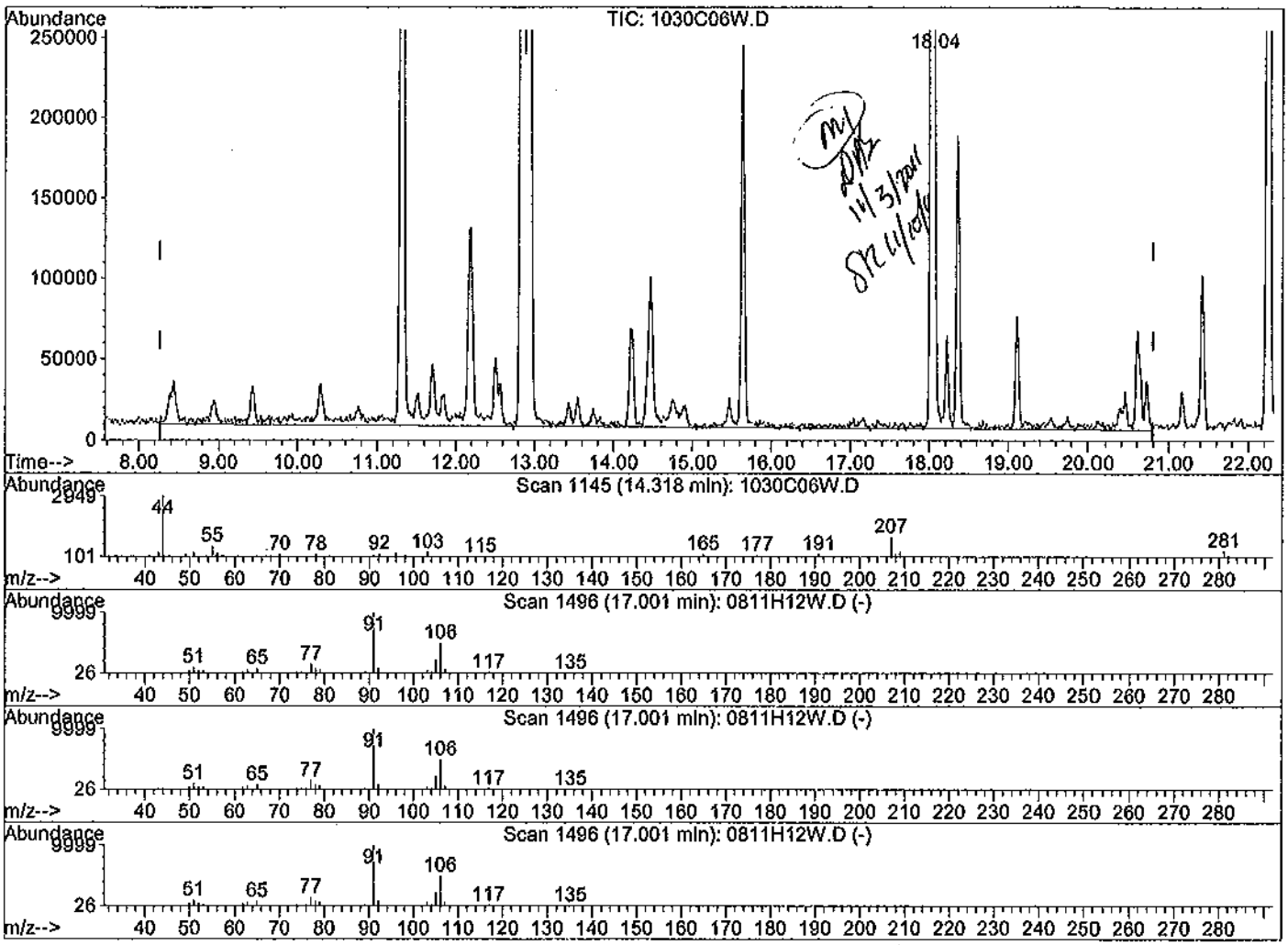
Ion	Exp%	Act%
TIC	100	100
0.00	0.00	0.68#
0.00	0.00	2.00#
0.00	0.00	0.00

1030C06W.D CGAS.M Thu Nov 03 10:30:25 2011

Quantitation Report

Data File : M:\CHICO\DATA\C111030\1030C06W.D Vial: 1
Acq On : 30 Oct 11 17:00 Operator: STC
Sample : Vol Std 10-30-11@50ug/L Inst : Chico
Misc : Water 10mLw/ IS:10-30-11 Multiplr: 1.00
Quant Time: Nov 3 10:30 2011 Quant Results File: temp.res

Method : M:\CHICO\DATA\C111030\CGAS.M (RTE Integrator)
Title : METHOD 8260B
Last Update : Mon Oct 31 09:35:31 2011
Response via : Multiple Level Calibration



TIC: 1030C06W.D

(2) Gasoline (TMHB)

18.04min 71.7166ppb m

response 17501250

Ion	Exp%	Act%
TIC	100	100
0.00	0.00	0.55#
0.00	0.00	1.62#
0.00	0.00	0.00

Quantitation Report (QT Reviewed)

Data File : M:\CHICO\DATA\C111030\1030C07W.D Vial: 1
Acq On : 30 Oct 11 17:43 Operator: STC
Sample : Vol Std 10-30-11@100ug/L Inst : Chico
Misc : Water 10mLw/ IS:10-30-11 Multiplr: 1.00

Quant Time: Nov 3 10:38 2011 Quant Results File: CGAS.RES

Quant Method : M:\CHICO\DATA\C111030\CGAS.M (RTE Integrator)
Title : METHOD 8260B
Last Update : Mon Oct 31 09:32:18 2011
Response via : Initial Calibration
DataAcq Meth : V8260

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Fluorobenzene (IS)	12.85	TIC	1049972	25.00000	ppb	0.00
3) Chlorobenzene-D5 (IS)	18.04	TIC	1057194	25.00000	ppb	0.00
4) 1,4-Dichlorobenzene-D (IS)	22.24	TIC	1054110	25.00000	ppb	0.00

System Monitoring Compounds

Target Compounds						Qvalue
2) Gasoline	18.04	TIC	21647604m	90.78273	ppb	100

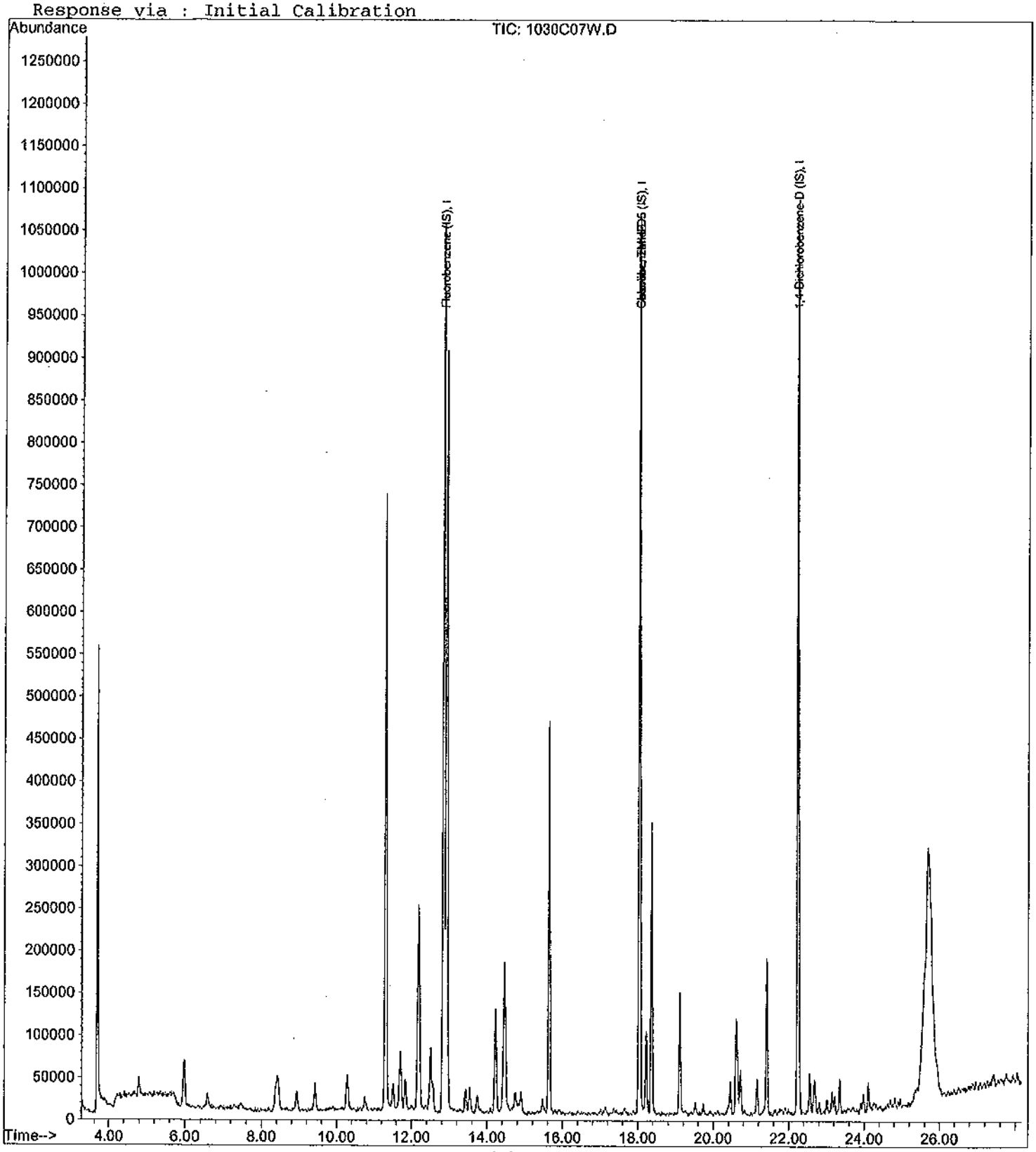
(#) = qualifier out of range (m) = manual integration
1030C07W.D CGAS.M Thu Nov 03 11:12:06 2011

Quantitation Report

Data File : M:\CHICO\DATA\C111030\1030C07W.D Vial: 1
Acq On : 30 Oct 11 17:43 Operator: STC
Sample : Vol Std 10-30-11@100ug/L Inst : Chico
Misc : Water 10mLw/ IS:10-30-11 Multiplr: 1.00

Quant Time: Nov 3 10:38 2011 Quant Results File: CGAS.RES

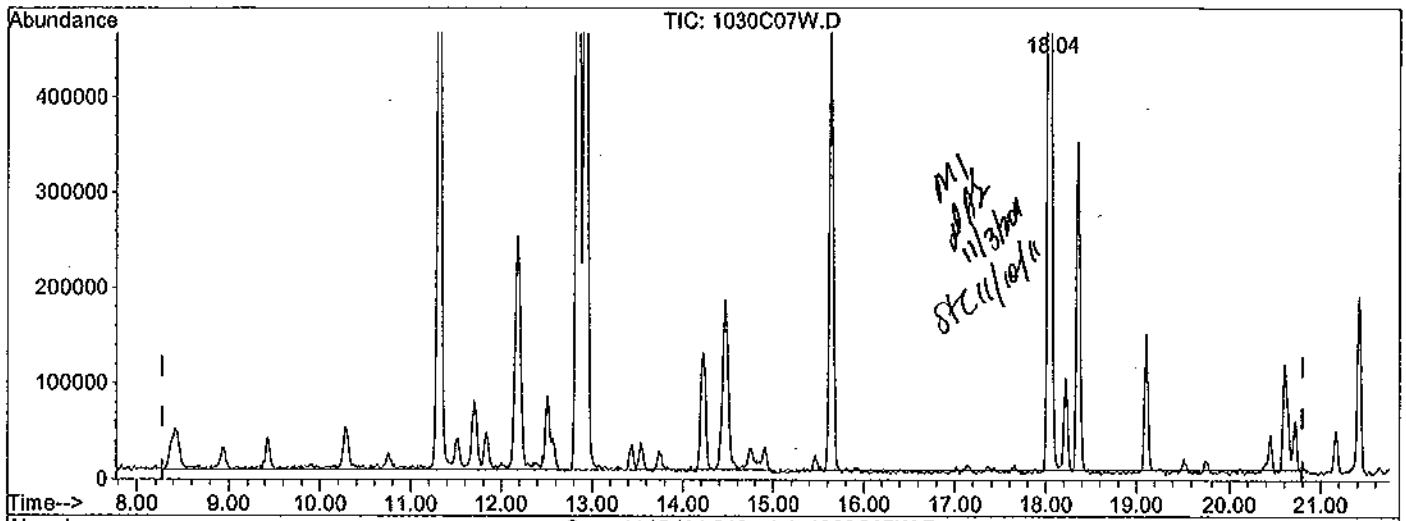
Method : M:\CHICO\DATA\C111030\CGAS.M (RTE Integrator)
Title : METHOD 8260B
Last Update : Thu Nov 03 10:47:02 2011
Response via : Initial Calibration

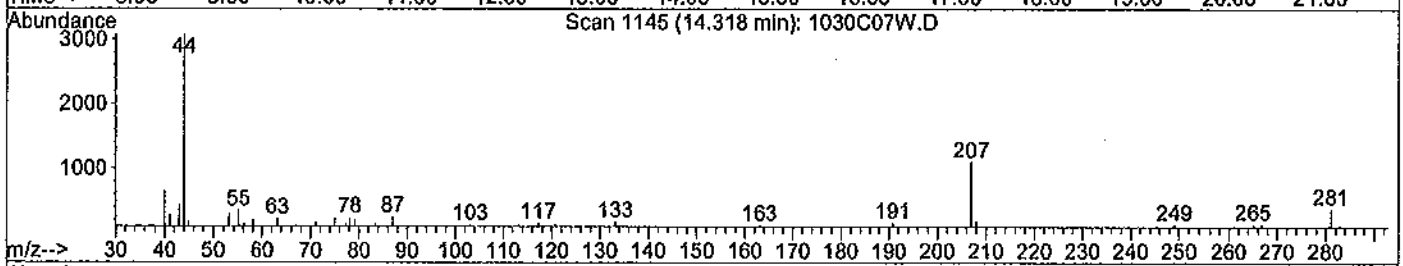


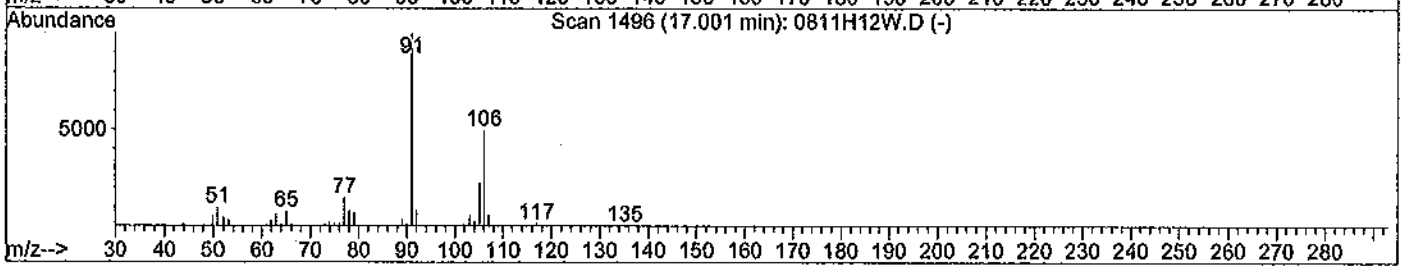
Quantitation Report

Data File : M:\CHICO\DATA\C111030\1030C07W.D Vial: 1
Acq On : 30 Oct 11 17:43 Operator: STC
Sample : Vol Std 10-30-11@100ug/L Inst : Chico
Misc : Water 10mLw/ IS:10-30-11 Multiplr: 1.00
Quant Time: Oct 31 9:32 2011 Quant Results File: temp.res

Method : M:\CHICO\DATA\C111030\CGAS.M (RTE Integrator)
Title : METHOD 8260B
Last Update : Mon Oct 31 09:35:31 2011
Response via : Multiple Level Calibration







TIC: 1030C07W.D

(2) Gasoline (TMHB)

14.31min 75.4746ppb m

response 17997299

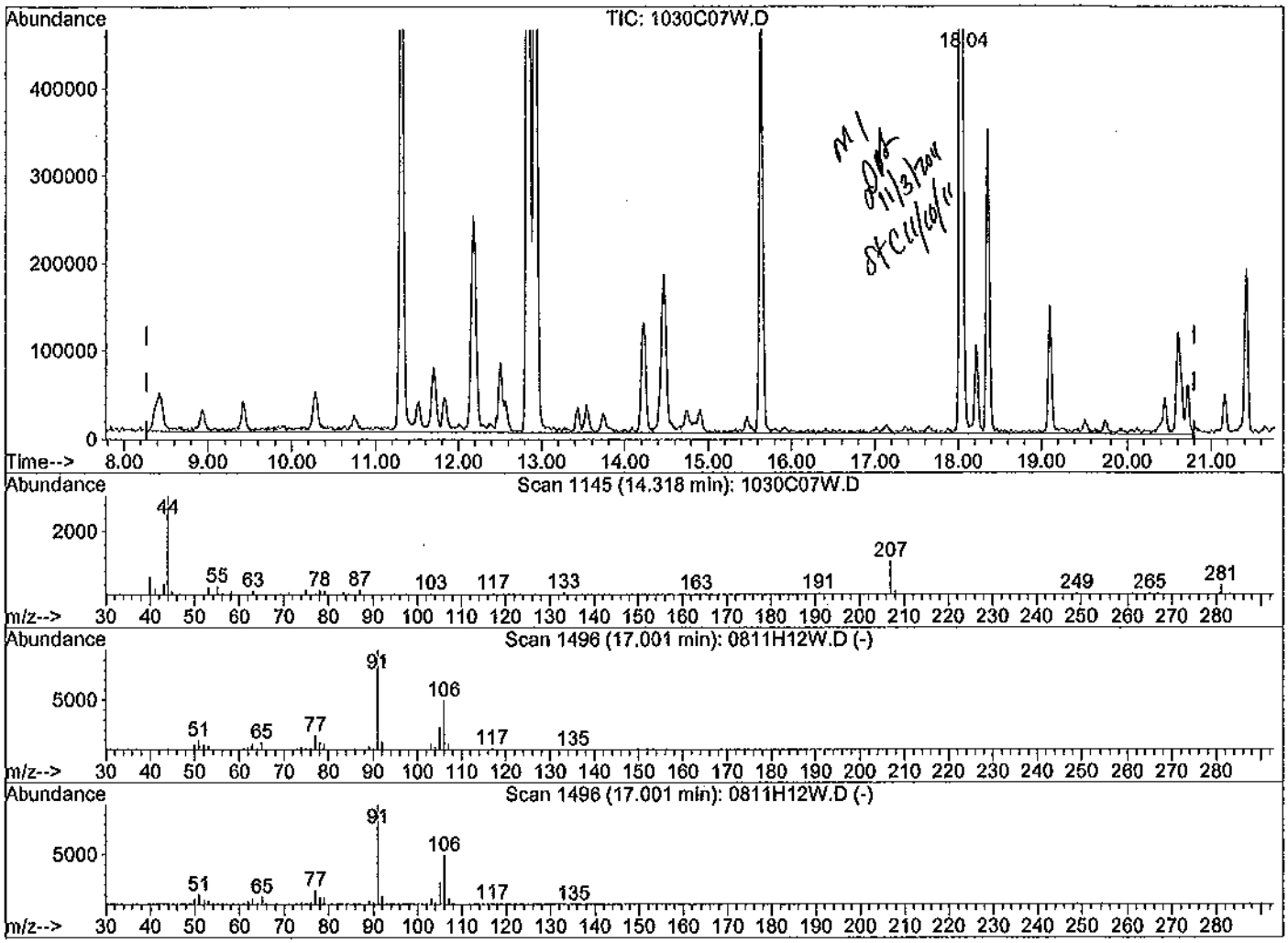
Ion	Exp%	Act%
TIC	100	100
0.00	0.00	0.55#
0.00	0.00	1.56#
0.00	0.00	0.00

1030C07W.D CGAS.M Thu Nov 03 10:31:29 2011

Quantitation Report

Data File : M:\CHICO\DATA\C111030\1030C07W.D Vial: 1
Acq On : 30 Oct 11 17:43 Operator: STC
Sample : Vol Std 10-30-11@100ug/L Inst : Chico
Misc : Water 10mLw/ IS:10-30-11 Multiplr: 1.00
Quant Time: Nov 3 10:38 2011 Quant Results File: temp.res

Method : M:\CHICO\DATA\C111030\CGAS.M (RTE Integrator)
Title : METHOD 8260B
Last Update : Mon Oct 31 09:35:31 2011
Response via : Multiple Level Calibration



TIC: 1030C07W.D

(2) Gasoline (TMHB)

18.04min 90.7827ppb m

response 21647604

Ion	Exp%	Act%
TIC	100	100
0.00	0.00	0.45#
0.00	0.00	1.30#
0.00	0.00	0.00

Quantitation Report (QT Reviewed)

Data File : M:\CHICO\DATA\C111030\1030C08W.D Vial: 1
Acq On : 30 Oct 11 18:26 Operator: STC
Sample : Vol Std 10-30-11@300ug/L Inst : Chico
Misc : Water 10mLw/ IS:10-30-11 Multiplr: 1.00

Quant Time: Nov 3 10:40 2011 Quant Results File: CGAS.RES

Quant Method : M:\CHICO\DATA\C111030\CGAS.M (RTE Integrator)
Title : METHOD 8260B
Last Update : Mon Oct 31 09:32:18 2011
Response via : Initial Calibration
DataAcq Meth : V8260

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Fluorobenzene (IS)	12.84	TIC	1085666	25.00000	ppb	0.00
3) Chlorobenzene-D5 (IS)	18.04	TIC	1080398	25.00000	ppb	0.00
4) 1,4-Dichlorobenzene-D (IS)	22.24	TIC	1118273	25.00000	ppb	0.00

System Monitoring Compounds

Target Compounds						Qvalue
2) Gasoline	15.64	TIC	39740510m	161.17894	ppb	100

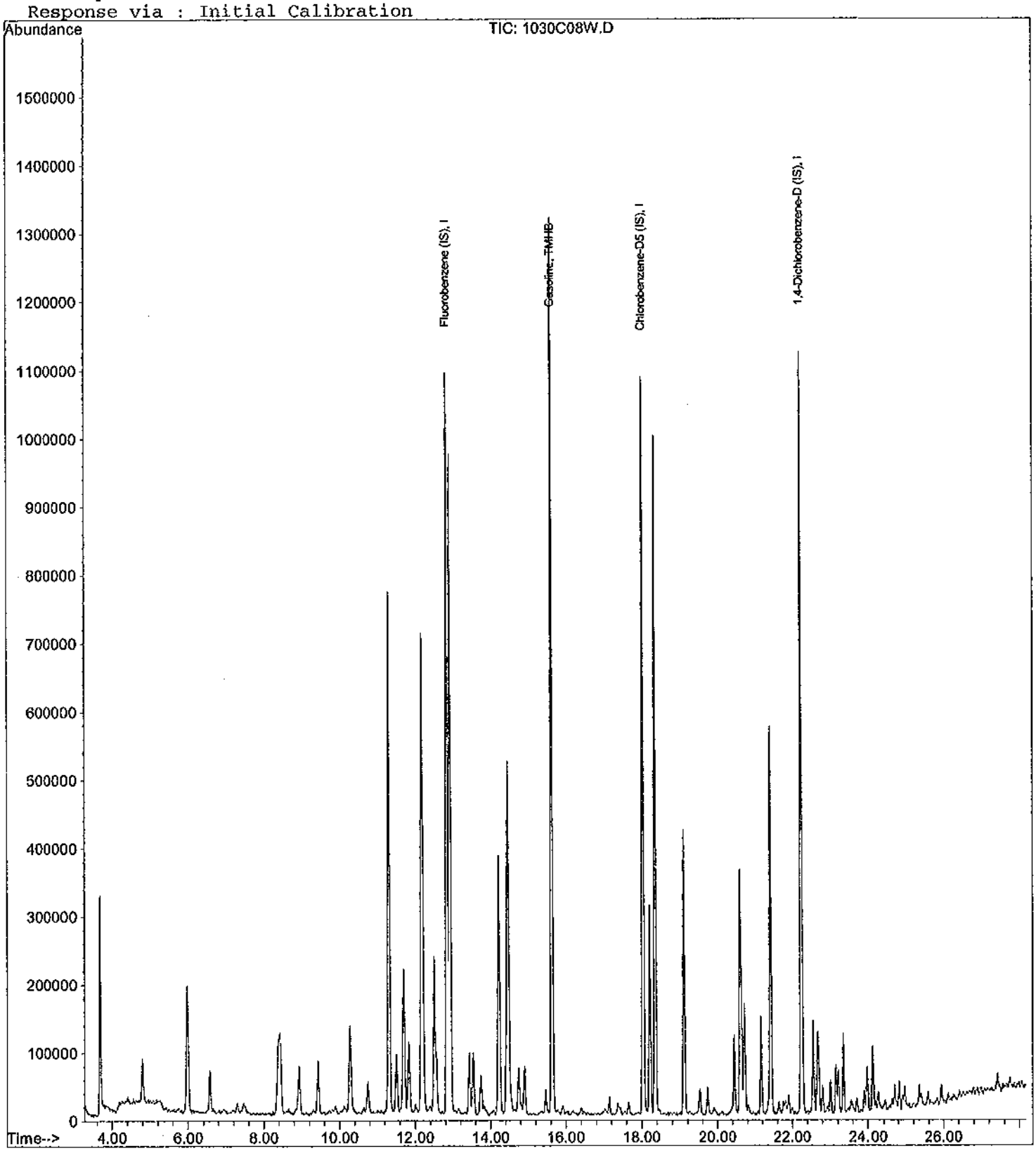
(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : M:\CHICO\DATA\C111030\1030C08W.D Vial: 1
Acq On : 30 Oct 11 18:26 Operator: STC
Sample : Vol Std 10-30-11@300ug/L Inst : Chico
Misc : Water 10mLw/ IS:10-30-11 Multiplr: 1.00

Quant Time: Nov 3 10:40 2011 Quant Results File: CGAS.RES

Method : M:\CHICO\DATA\C111030\CGAS.M (RTE Integrator)
Title : METHOD 8260B
Last Update : Thu Nov 03 10:47:02 2011
Response via : Initial Calibration



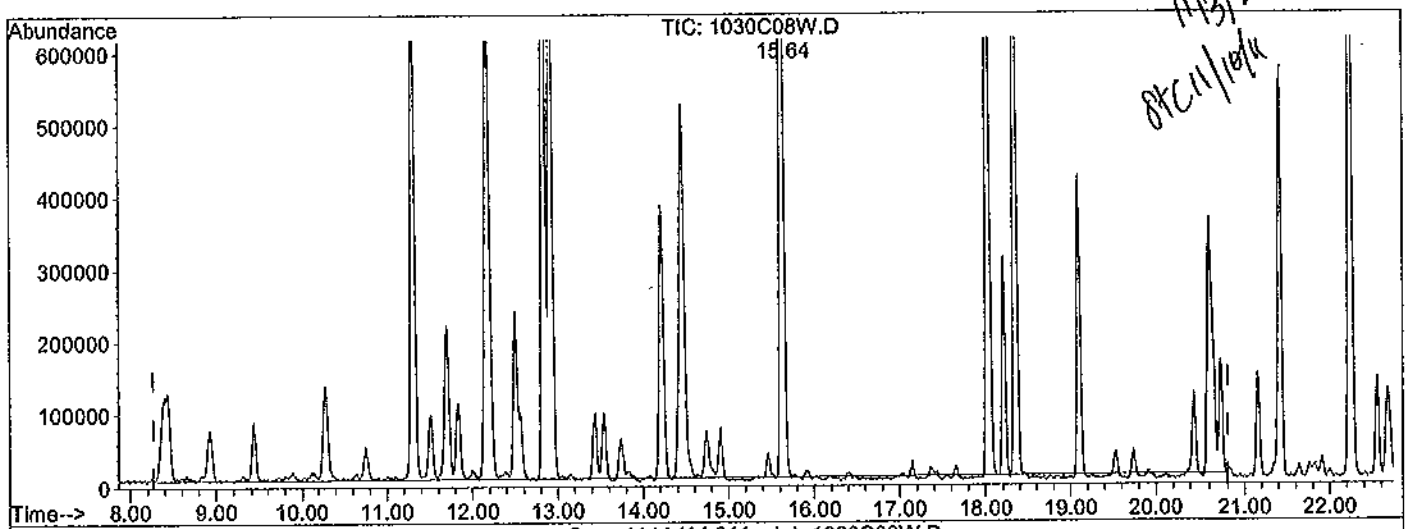
Quantitation Report

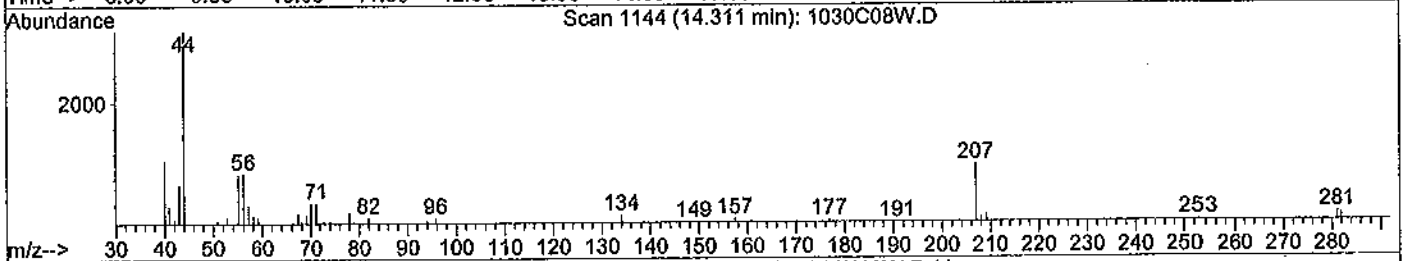
Data File : M:\CHICO\DATA\C111030\1030C08W.D
Acq On : 30 Oct 11 18:26
Sample : Vol Std 10-30-11@300ug/L
Misc : Water 10mLw/ IS:10-30-11
Quant Time: Oct 31 9:32 2011

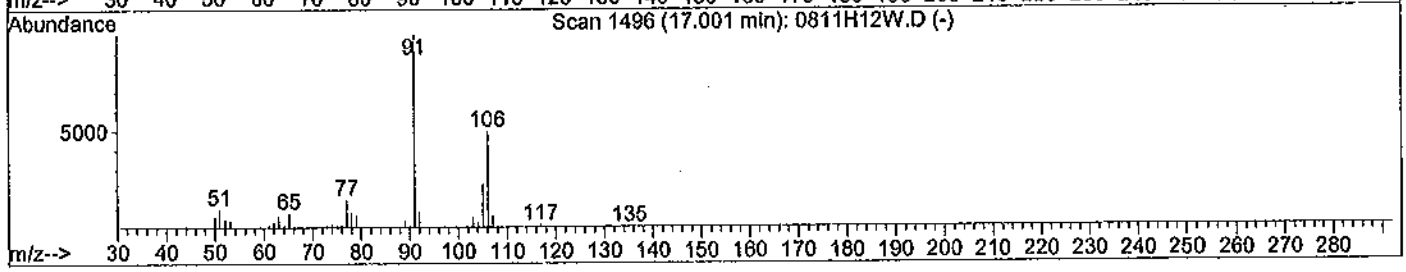
Vial: 1
Operator: STC
Inst : Chico
Multiplr: 1.00
Quant Results File: temp.res

Method : M:\CHICO\DATA\C111030\CGAS.M (RTE Integrator)
Title : METHOD 8260B
Last Update : Mon Oct 31 09:35:31 2011
Response via : Multiple Level Calibration

11/3/2011 STC







TIC: 1030C08W.D

(2) Gasoline (TMHB)

14.31min 137.6327ppb m

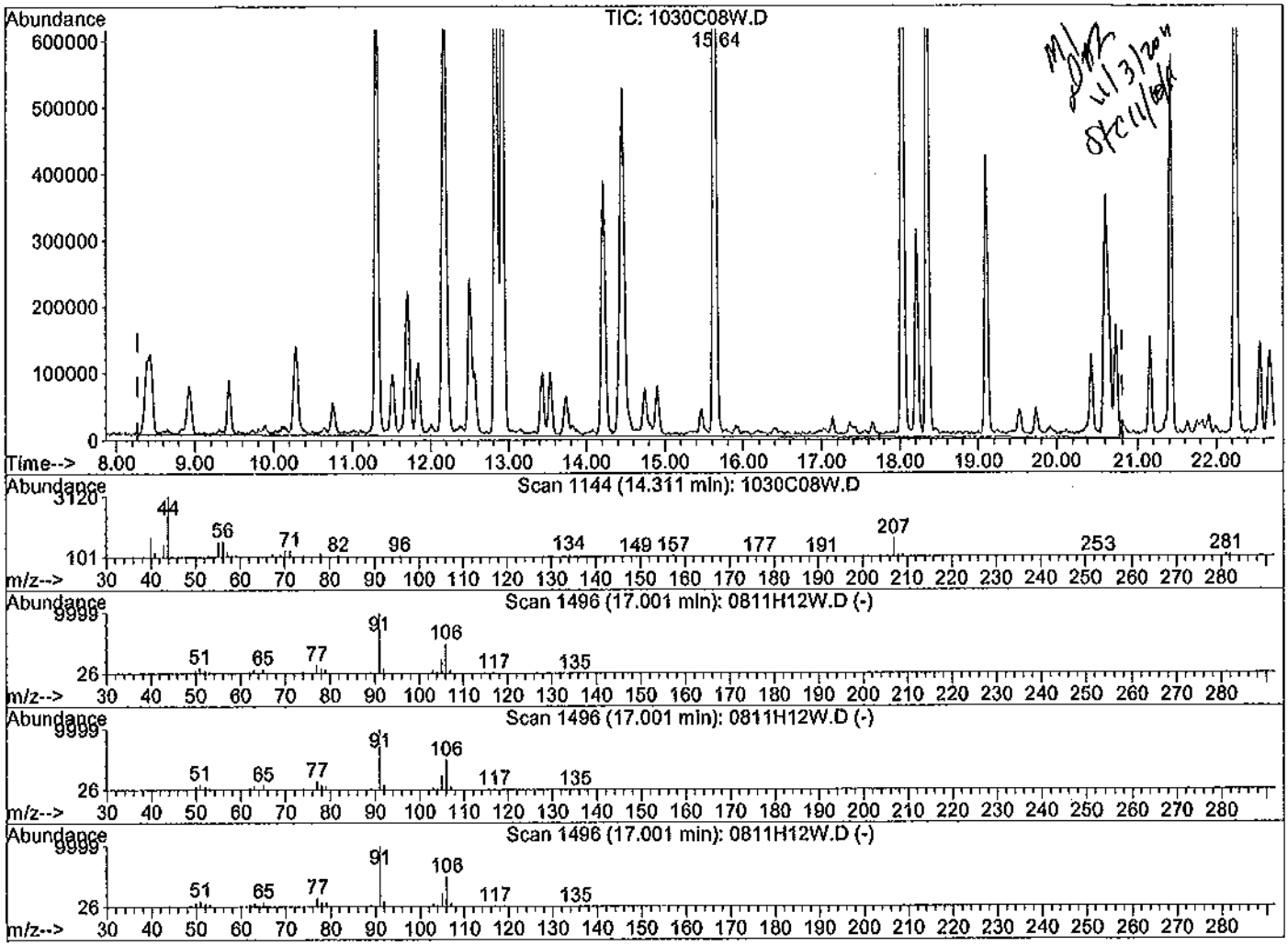
response 33934923

Ion	Exp%	Act%
TIC	100	100
0.00	0.00	0.30#
0.00	0.00	0.88#
0.00	0.00	0.00

Quantitation Report

Data File : M:\CHICO\DATA\C111030\1030C08W.D Vial: 1
Acq On : 30 Oct 11 18:26 Operator: STC
Sample : Vol Std 10-30-11@300ug/L Inst : Chico
Misc : Water 10mLw/ IS:10-30-11 Multiplr: 1.00
Quant Time: Nov 3 10:40 2011 Quant Results File: temp.res

Method : M:\CHICO\DATA\C111030\CGAS.M (RTE Integrator)
Title : METHOD 8260B
Last Update : Mon Oct 31 09:35:31 2011
Response via : Multiple Level Calibration



TIC: 1030C08W.D

(2) Gasoline (TMHB)

15.64min 161.1789ppb m

response 39740510

Ion	Exp%	Act%
TIC	100	100
0.00	0.00	0.25#
0.00	0.00	0.75#
0.00	0.00	0.00

Quantitation Report (QT Reviewed)

Data File : M:\CHICO\DATA\C111030\1030C09W.D Vial: 1
Acq On : 30 Oct 11 19:09 Operator: STC
Sample : Vol Std 10-30-11@600ug/L Inst : Chico
Misc : Water 10mLw/ IS:10-30-11 Multiplr: 1.00

Quant Time: Nov 3 10:41 2011 Quant Results File: CGAS.RES

Quant Method : M:\CHICO\DATA\C111030\CGAS.M (RTE Integrator)
Title : METHOD 8260B
Last Update : Mon Oct 31 09:32:18 2011
Response via : Initial Calibration
DataAcq Meth : V8260

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Fluorobenzene (IS)	12.84	TIC	1104080	25.00000	ppb	0.00
3) Chlorobenzene-D5 (IS)	18.04	TIC	1114811	25.00000	ppb	0.00
4) 1,4-Dichlorobenzene-D (IS)	22.25	TIC	1175050	25.00000	ppb	0.00

System Monitoring Compounds

Target Compounds						Qvalue
2) Gasoline	15.64	TIC	65808275m	262.45271	ppb	100

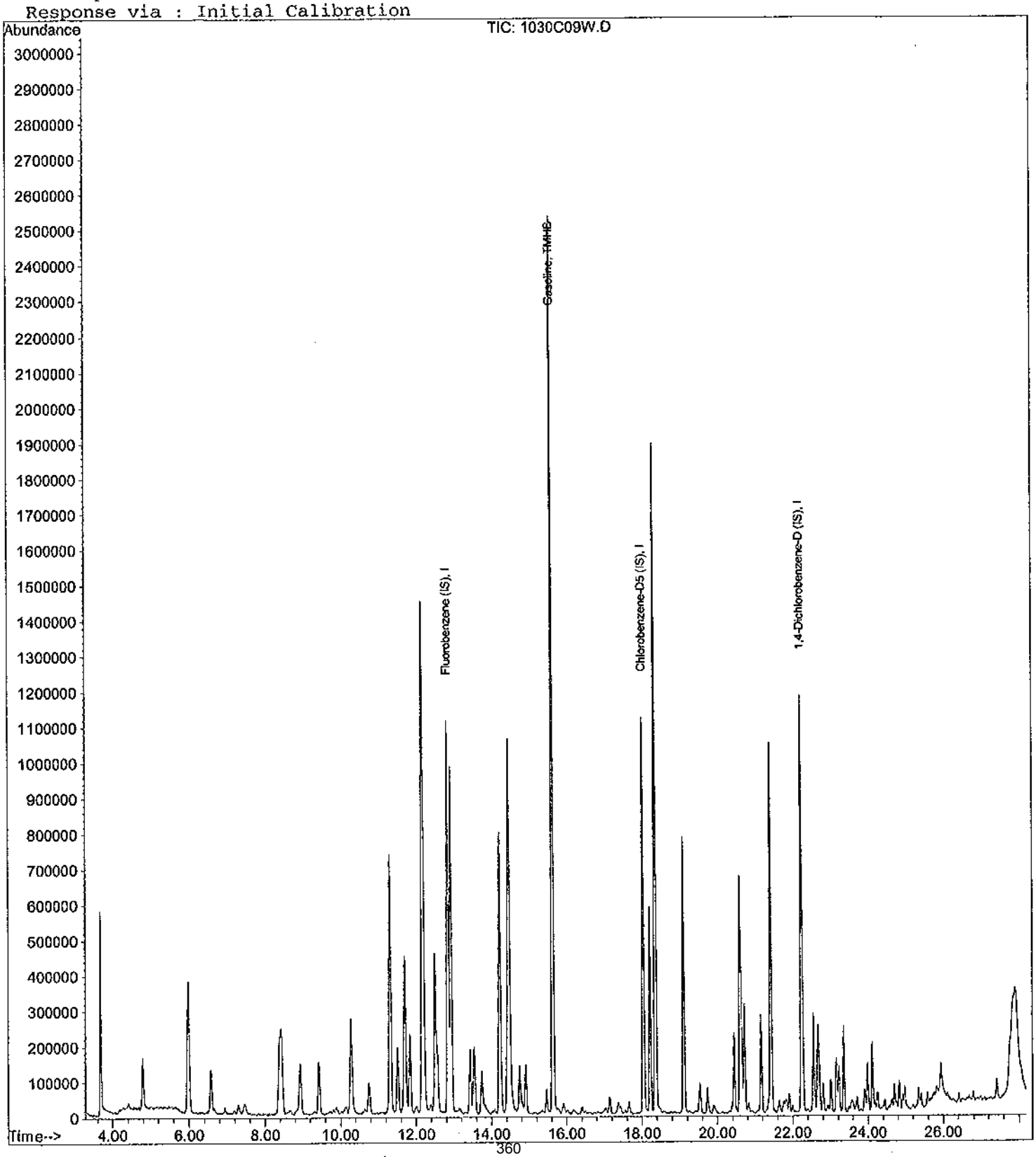
(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : M:\CHICO\DATA\C111030\1030C09W.D Vial: 1
Acq On : 30 Oct 11 19:09 Operator: STC
Sample : Vol Std 10-30-11@600ug/L Inst : Chico
Misc : Water 10mLw/ IS:10-30-11 Multiplr: 1.00

Quant Time: Nov 3 10:41 2011 Quant Results File: CGAS.RES

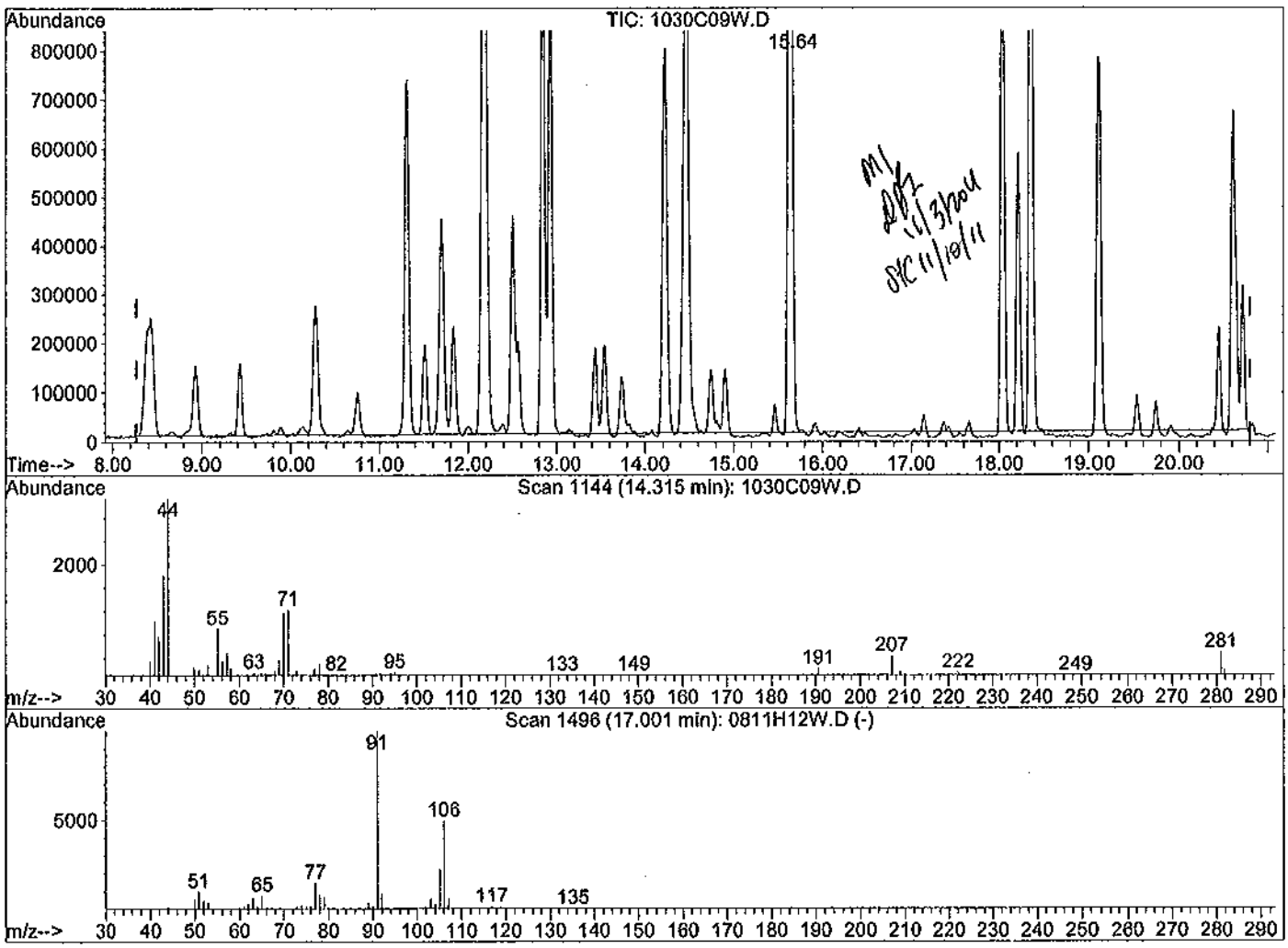
Method : M:\CHICO\DATA\C111030\CGAS.M (RTE Integrator)
Title : METHOD 8260B
Last Update : Thu Nov 03 10:47:02 2011
Response via : Initial Calibration



Quantitation Report

Data File : M:\CHICO\DATA\C111030\1030C09W.D Vial: 1
Acq On : 30 Oct 11 19:09 Operator: STC
Sample : Vol Std 10-30-11@600ug/L Inst : Chico
Misc : Water 10mLw/ IS:10-30-11 Multiplr: 1.00
Quant Time: Oct 31 9:32 2011 Quant Results File: temp.res

Method : M:\CHICO\DATA\C111030\CGAS.M (RTE Integrator)
Title : METHOD 8260B
Last Update : Mon Oct 31 09:35:31 2011
Response via : Multiple Level Calibration



TIC: 1030C09W.D

(2) Gasoline (TMHB)

14.31min 231.1564ppb m

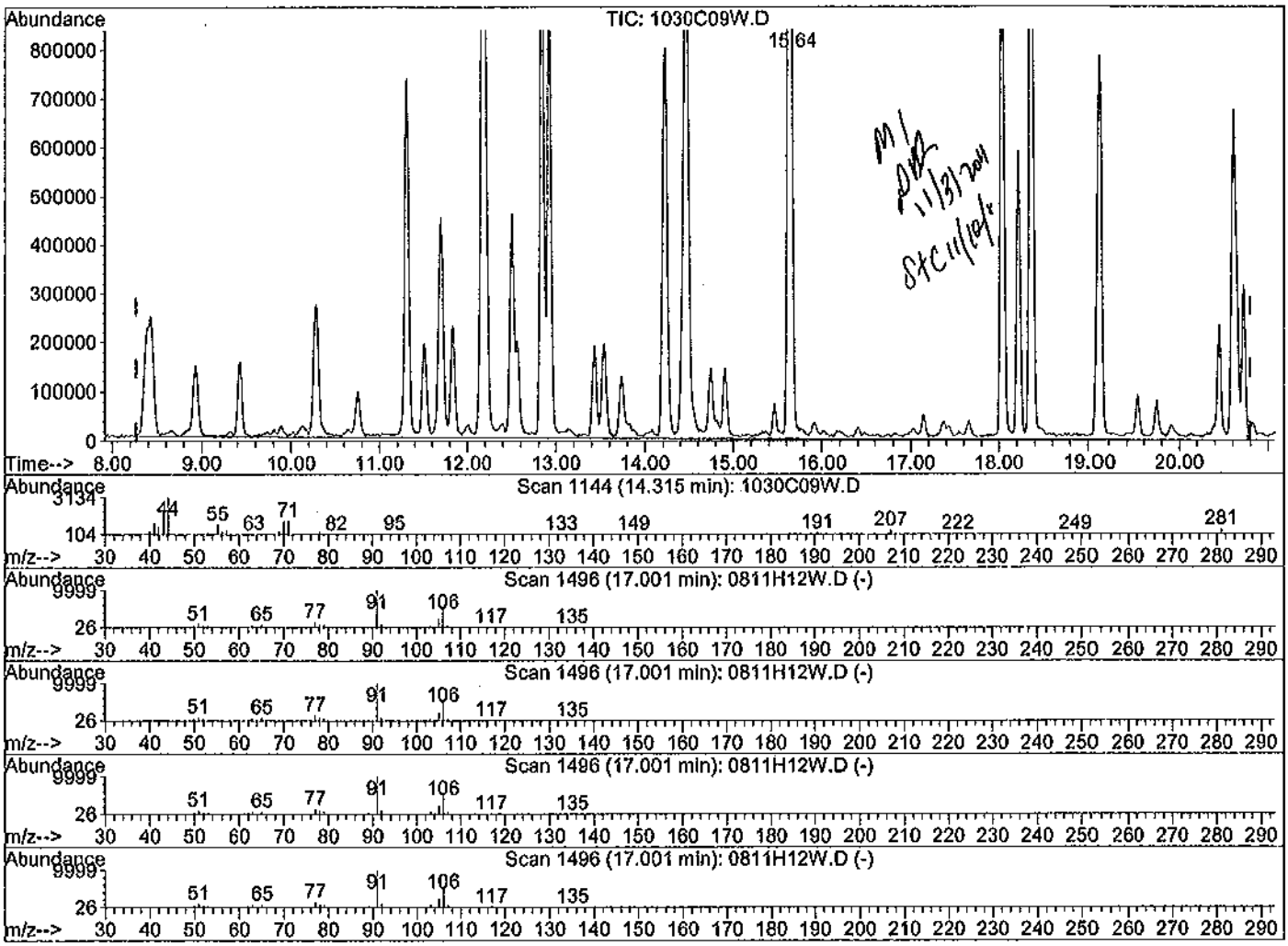
response 57960938

Ion	Exp%	Act%
TIC	100	100
0.00	0.00	0.18#
0.00	0.00	0.51#
0.00	0.00	0.00

Quantitation Report

Data File : M:\CHICO\DATA\C111030\1030C09W.D Vial: 1
Acq On : 30 Oct 11 19:09 Operator: STC
Sample : Vol Std 10-30-11@600ug/L Inst : Chico
Misc : Water 10mLw/ IS:10-30-11 Multiplr: 1.00
Quant Time: Nov 3 10:41 2011 Quant Results File: temp.res

Method : M:\CHICO\DATA\C111030\CGAS.M (RTE Integrator)
Title : METHOD 8260B
Last Update : Mon Oct 31 09:35:31 2011
Response via : Multiple Level Calibration



TIC: 1030C09W.D

(2) Gasoline (TMHB)

15.64min 262.4527ppb m

response 65808275

Ion	Exp%	Act%
TIC	100	100
0.00	0.00	0.16#
0.00	0.00	0.45#
0.00	0.00	0.00

Quantitation Report (QT Reviewed)

Data File : M:\CHICO\DATA\C111030\1030C10W.D Vial: 1
Acq On : 30 Oct 11 19:52 Operator: STC
Sample : Vol Std 10-30-11@800ug/L Inst : Chico
Misc : Water 10mLw/ IS:10-30-11 Multiplr: 1.00

Quant Time: Nov 3 10:42 2011 Quant Results File: CGAS.RES

Quant Method : M:\CHICO\DATA\C111030\CGAS.M (RTE Integrator)
Title : METHOD 8260B
Last Update : Mon Oct 31 09:32:18 2011
Response via : Initial Calibration
DataAcq Meth : V8260

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Fluorobenzene (IS)	12.84	TIC	1129347	25.00000	ppb	0.00
3) Chlorobenzene-D5 (IS)	18.03	TIC	1159453	25.00000	ppb	0.00
4) 1,4-Dichlorobenzene-D (IS)	22.24	TIC	1268278	25.00000	ppb	0.00

System Monitoring Compounds

Target Compounds						Qvalue
2) Gasoline	15.63	TIC	84666447m	330.10723	ppb	100

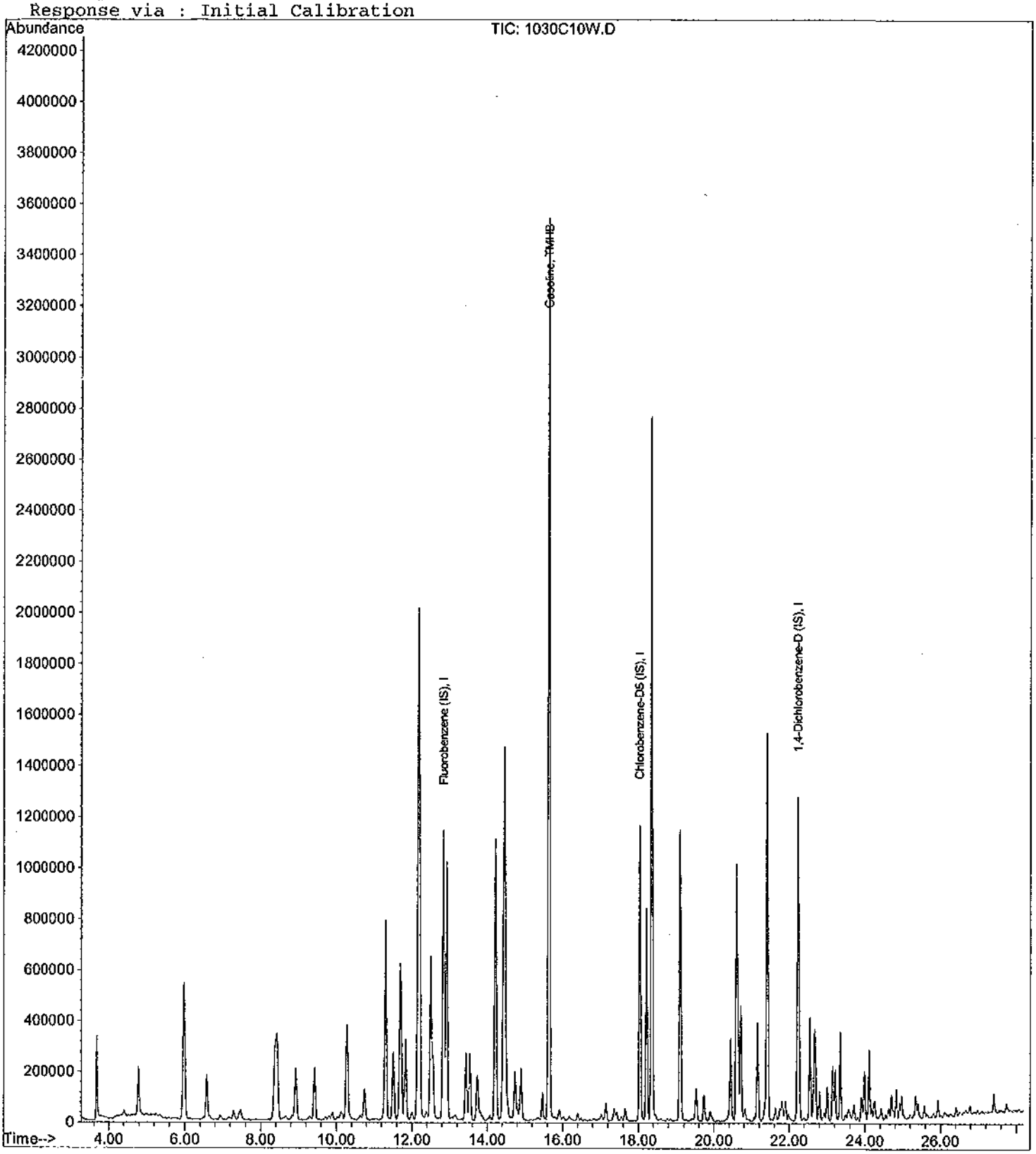
(#) = qualifier out of range (m) = manual integration
1030C10W.D CGAS.M Thu Nov 03 11:12:19 2011 Page 1

Quantitation Report

Data File : M:\CHICO\DATA\C111030\1030C10W.D Vial: 1
Acq On : 30 Oct 11 19:52 Operator: STC
Sample : Vol Std 10-30-11@800ug/L Inst : Chico
Misc : Water 10mLw/ IS:10-30-11 Multiplr: 1.00

Quant Time: Nov 3 10:42 2011 Quant Results File: CGAS.RES

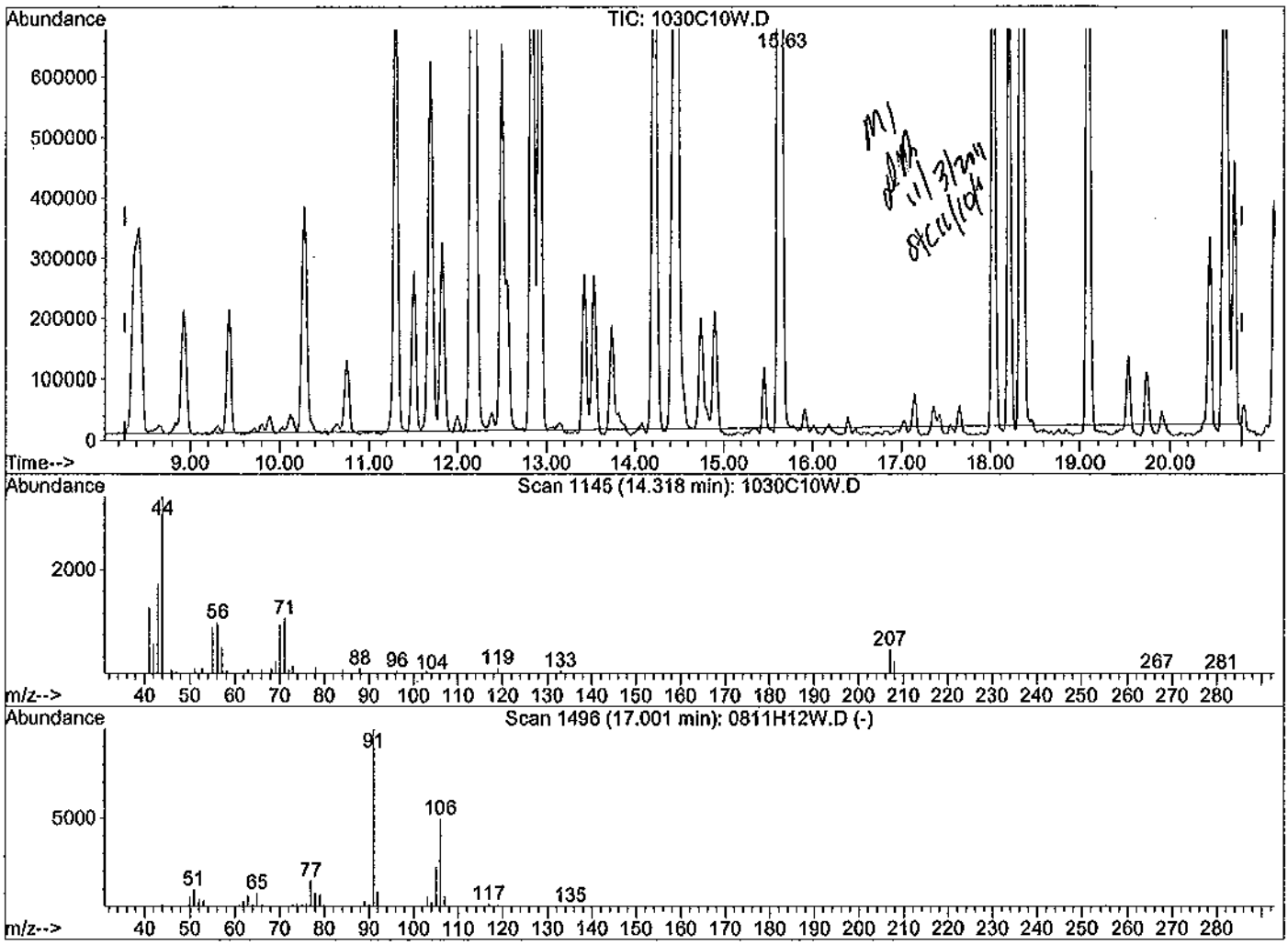
Method : M:\CHICO\DATA\C111030\CGAS.M (RTE Integrator)
Title : METHOD 8260B
Last Update : Thu Nov 03 10:47:02 2011
Response via : Initial Calibration



Quantitation Report

Data File : M:\CHICO\DATA\C111030\1030C10W.D Vial: 1
Acq On : 30 Oct 11 19:52 Operator: STC
Sample : Vol Std 10-30-11@800ug/L Inst : Chico
Misc : Water 10mLw/ IS:10-30-11 Multiplr: 1.00
Quant Time: Oct 31 9:32 2011 Quant Results File: temp.res

Method : M:\CHICO\DATA\C111030\CGAS.M (RTE Integrator)
Title : METHOD 8260B
Last Update : Mon Oct 31 09:35:31 2011
Response via : Multiple Level Calibration



TIC: 1030C10W.D

(2) Gasoline (TMHB)

14.31min 303.9125ppb m

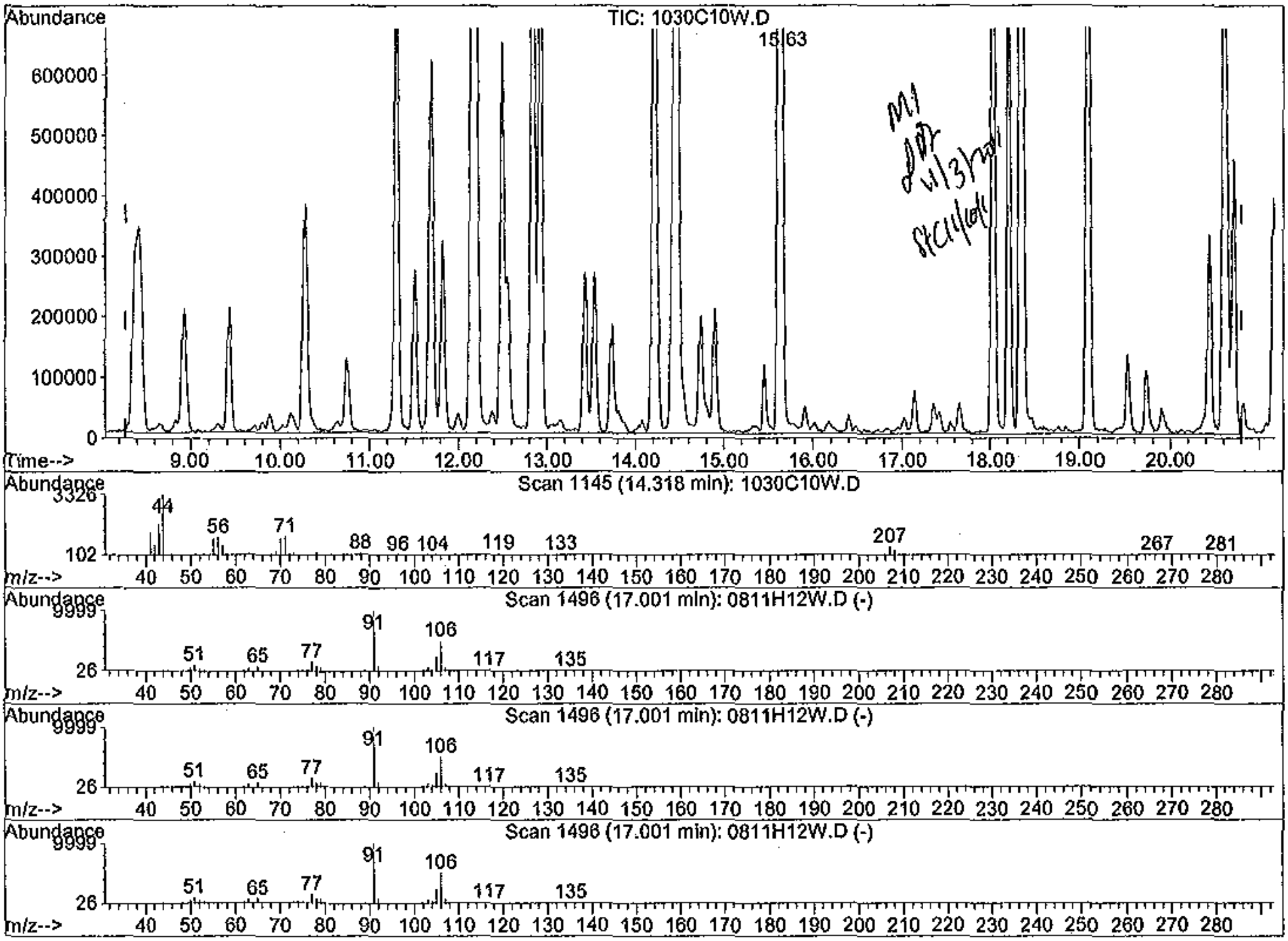
response 77947975

Ion	Exp%	Act%
TIC	100	100
0.00	0.00	0.13#
0.00	0.00	0.41#
0.00	0.00	0.00

Quantitation Report

```
Data File : M:\CHICO\DATA\C111030\1030C10W.D              Vial: 1
Acq On    : 30 Oct 11  19:52                          Operator: STC
Sample    : Vol Std 10-30-11@800ug/L                  Inst    : Chico
Misc      : Water 10mLw/ IS:10-30-11                  Multiplr: 1.00
Quant Time: Nov  3 10:42 2011               Quant Results File: temp.res

Method        : M:\CHICO\DATA\C111030\CGAS.M (RTE Integrator)
Title         : METHOD 8260B
Last Update   : Mon Oct 31 09:35:31 2011
Response via  : Multiple Level Calibration
```



TIC: 1030C10W.D

(2) Gasoline (TMHB)

15.63min 330.1072ppb m

response 84666447

Ion	Exp%	Act%
TIC	100	100
0.00	0.00	0.12#
0.00	0.00	0.38#
0.00	0.00	0.00

1030C10W.D CGAS.M Thu Nov 03 10:42:47 2011

Quantitation Report (QT Reviewed)

Data File : M:\CHICO\DATA\C111030\1030C11W.D Vial: 1
Acq On : 30 Oct 11 20:35 Operator: STC
Sample : Vol Std 10-30-i1@1000ug/L Inst : Chico
Misc : Water 10mLw/ IS:10-30-11 Multiplr: 1.00

Quant Time: Nov 3 10:43 2011 Quant Results File: CGAS.RES

Quant Method : M:\CHICO\DATA\C111030\CGAS.M (RTE Integrator)
Title : METHOD 8260B
Last Update : Mon Oct 31 09:32:18 2011
Response via : Initial Calibration
DataAcq Meth : V8260

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Fluorobenzene (IS)	12.84	TIC	1162372	25.00000	ppb	0.00
3) Chlorobenzene-D5 (IS)	18.03	TIC	1207961	25.00000	ppb	0.00
4) 1,4-Dichlorobenzene-D (IS)	22.24	TIC	1354742	25.00000	ppb	0.00

System Monitoring Compounds

Target Compounds						Qvalue
2) Gasoline	15.63	TIC	105748641m	400.59060	ppb	100

(#) = qualifier out of range (m) = manual integration

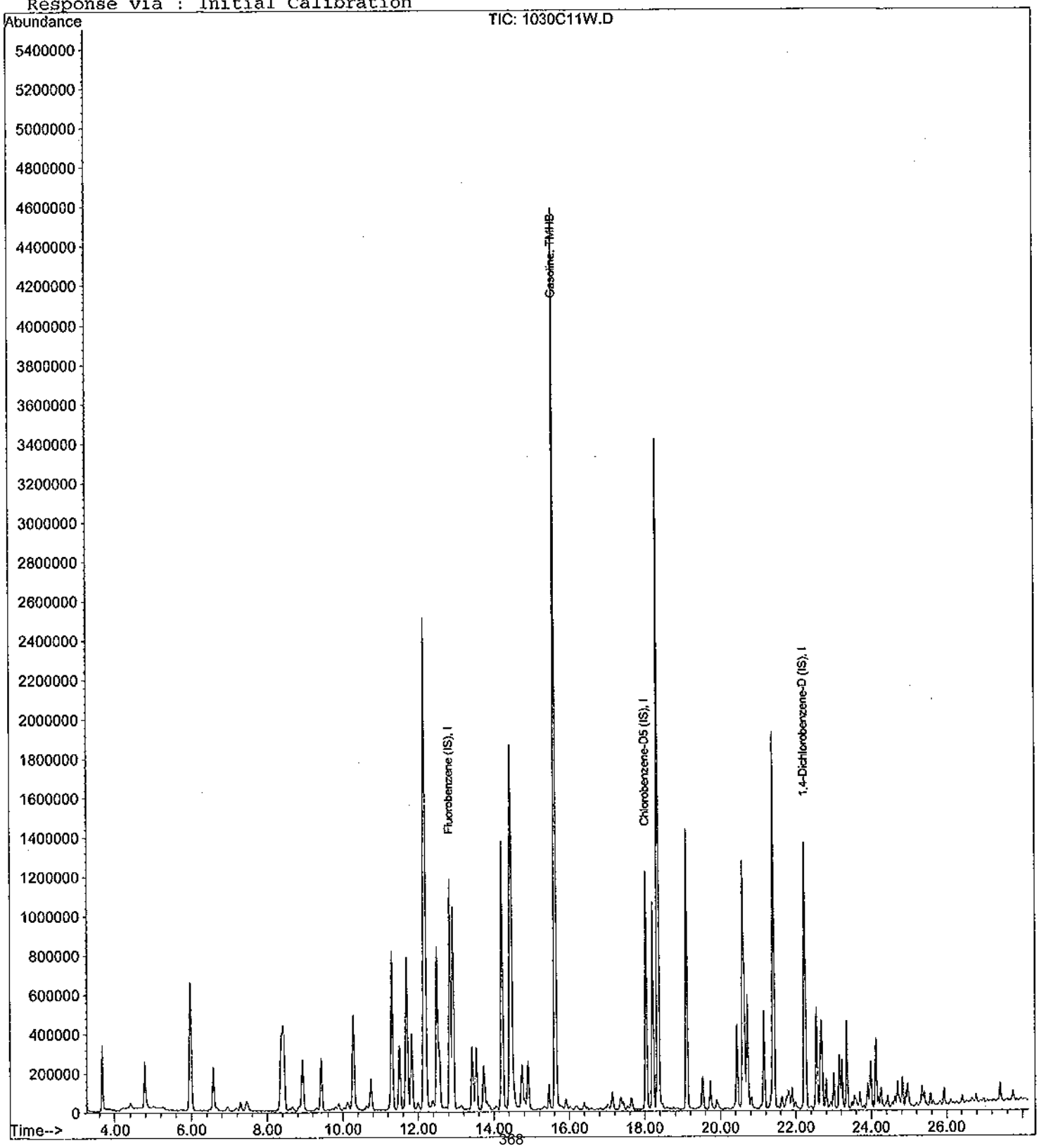
Quantitation Report

Data File : M:\CHICO\DATA\C111030\1030C11W.D
Acq On : 30 Oct 11 20:35
Sample : Vol Std 10-30-11@1000ug/L
Misc : Water 10mLw/ IS:10-30-11

Vial: 1
Operator: STC
Inst : Chico
Multiplr: 1.00

Quant Time: Nov 3 10:43 2011 Quant Results File: CGAS.RES

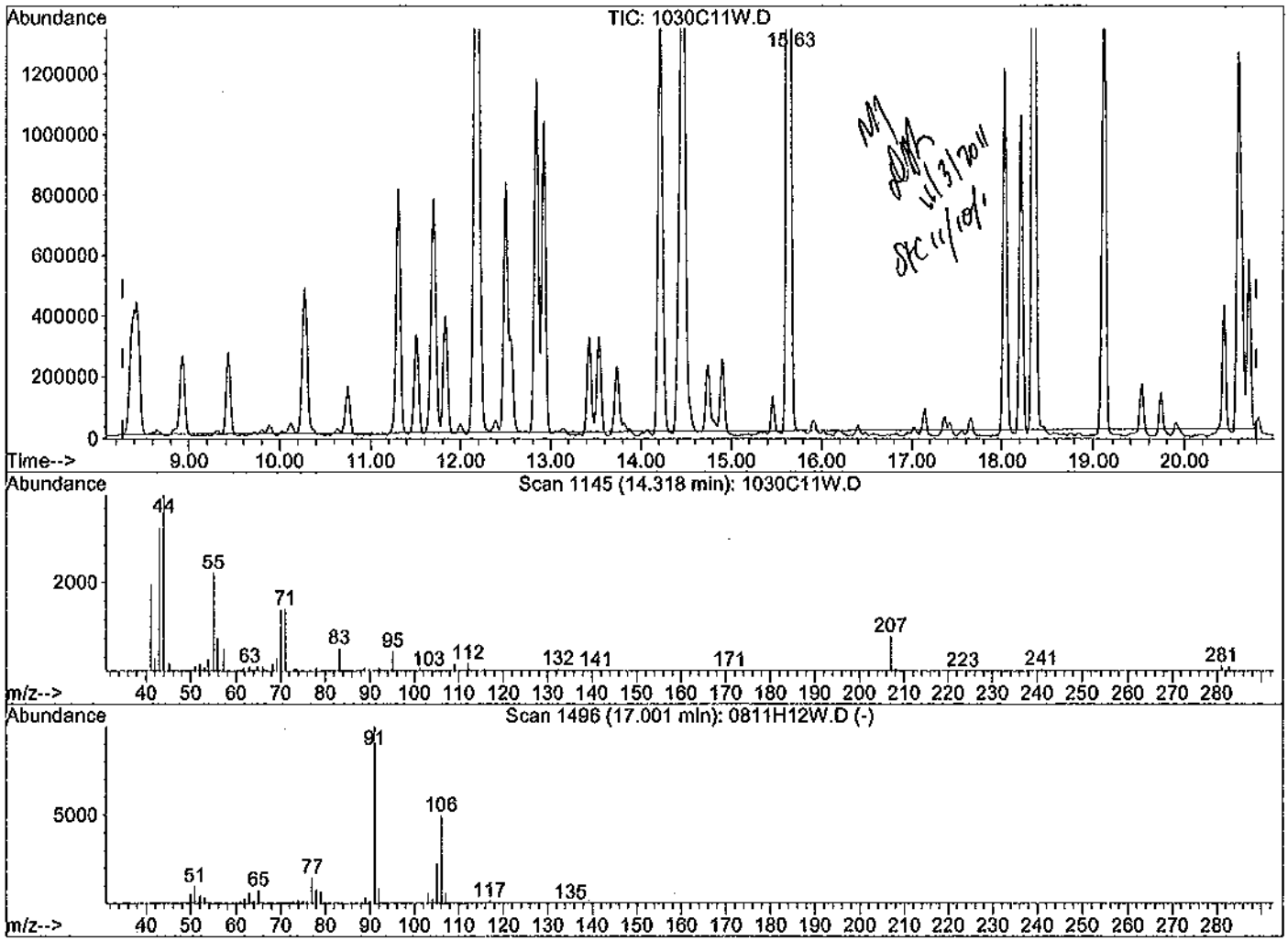
Method : M:\CHICO\DATA\C111030\CGAS.M (RTE Integrator)
Title : METHOD 8260B
Last Update : Thu Nov 03 10:47:02 2011
Response via : Initial Calibration



Quantitation Report

Data File : M:\CHICO\DATA\C111030\1030C11W.D Vial: 1
Acq On : 30 Oct 11 20:35 Operator: STC
Sample : Vol Std 10-30-11@1000ug/L Inst : Chico
Misc : Water 10mLw/ IS:10-30-11 Multiplr: 1.00
Quant Time: Oct 31 9:33 2011 Quant Results File: temp.res

Method : M:\CHICO\DATA\C111030\CGAS.M (RTE Integrator)
Title : METHOD 8260B
Last Update : Mon Oct 31 09:35:31 2011
Response via : Multiple Level Calibration



TIC: 1030C11W.D

(2) Gasoline (TMHB)

14.31min 368.6230ppb m

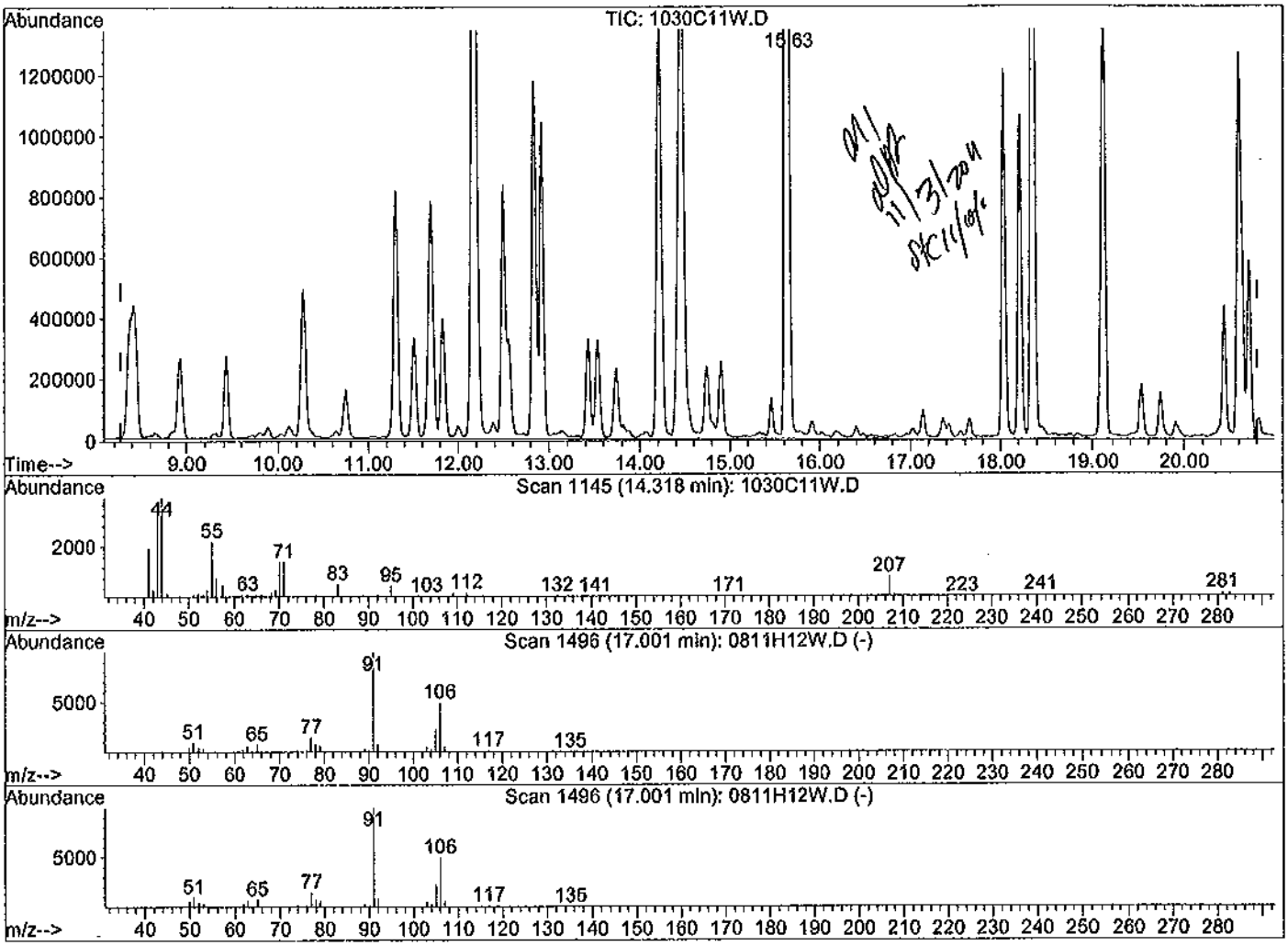
response 97309775

Ion	Exp%	Act%
TIC	100	100
0.00	0.00	0.11#
0.00	0.00	0.34#
0.00	0.00	0.00

Quantitation Report

Data File : M:\CHICO\DATA\C111030\1030C11W.D Vial: 1
Acq On : 30 Oct 11 20:35 Operator: STC
Sample : Vol Std 10-30-11@1000ug/L Inst : Chico
Misc : Water 10mLw/ IS:10-30-11 Multiplr: 1.00
Quant Time: Nov 3 10:43 2011 Quant Results File: temp.res

Method : M:\CHICO\DATA\C111030\CGAS.M (RTE Integrator)
Title : METHOD 8260B
Last Update : Mon Oct 31 09:35:31 2011
Response via : Multiple Level Calibration



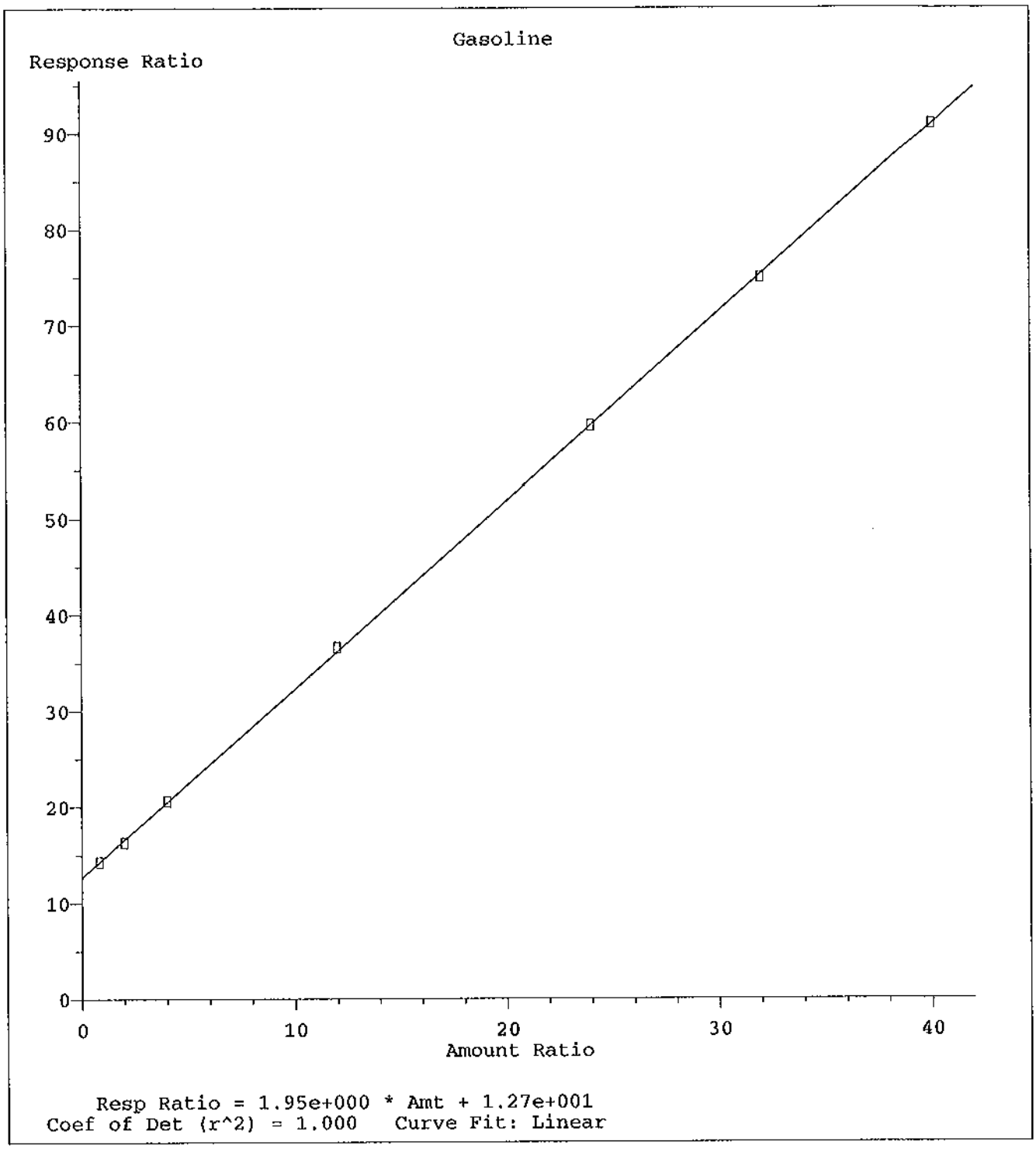
TIC: 1030C11W.D

(2) Gasoline (TMHB)

15.63min 400.5906ppb m

response 105748641

Ion	Exp%	Act%
TIC	100	100
0.00	0.00	0.10#
0.00	0.00	0.31#
0.00	0.00	0.00



Method Name: M:\CHICO\DATA\C111030\CGAS.M
Calibration Table Last Updated: Thu Nov 03 10:47:02 2011

VOLATILE ORGANIC ANALYSIS BY
EPA METHOD 8260B

Form 7

Second Source Calibration

Lab Name: APPL, Inc.
Case No:
Matrix:

SDG No:
Date Analyzed: 10/31/2011
Instrument: Chico
Initial Cal. Date: 10/30/2011
Data File: 1030C29W.D

		Compound	MEAN	CCRF	%D		%Drift
1	I	Fluorobenzene (IS)	ISTD			I	
2	TMHB	Gasoline	5.897	3.226	45	TMHBL	11
3	I	Chlorobenzene-D5 (IS)	ISTD			I	
4	I	1,4-Dichlorobenzene-D (IS)	ISTD			I	
5							
6							
7							
8							
9							
10							
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33							
34							
35							
36							
37							
38							
39							
40							
		Average			45.0		

Quantitation Report (QT Reviewed)

Data File : M:\CHICO\DATA\C111030\1030C29W.D Vial: 1
Acq On : 31 Oct 11 9:31 Operator: STC
Sample : GAS 300ug/L (SS) Inst : Chico
Misc : Water 10mLw/ IS&S:10-30/10-26-11 Multiplr: 1.00

Quant Time: Nov 3 10:51 2011 Quant Results File: CGAS.RES

Quant Method : M:\CHICO\DATA\C111030\CGAS.M (RTE Integrator)
Title : METHOD 8260B
Last Update : Thu Nov 03 10:47:02 2011
Response via : Initial Calibration
DataAcq Meth : V8260

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Fluorobenzene (IS)	12.84	TIC	1211423	25.00000	ppb	0.00
3) Chlorobenzene-D5 (IS)	18.03	TIC	1191079	25.00000	ppb	0.00
4) 1,4-Dichlorobenzene-D (IS)	22.24	TIC	1217266	25.00000	ppb	0.00

System Monitoring Compounds

Target Compounds						Qvalue
2) Gasoline	15.63	TIC	46900368m	332.66187	ppb	100

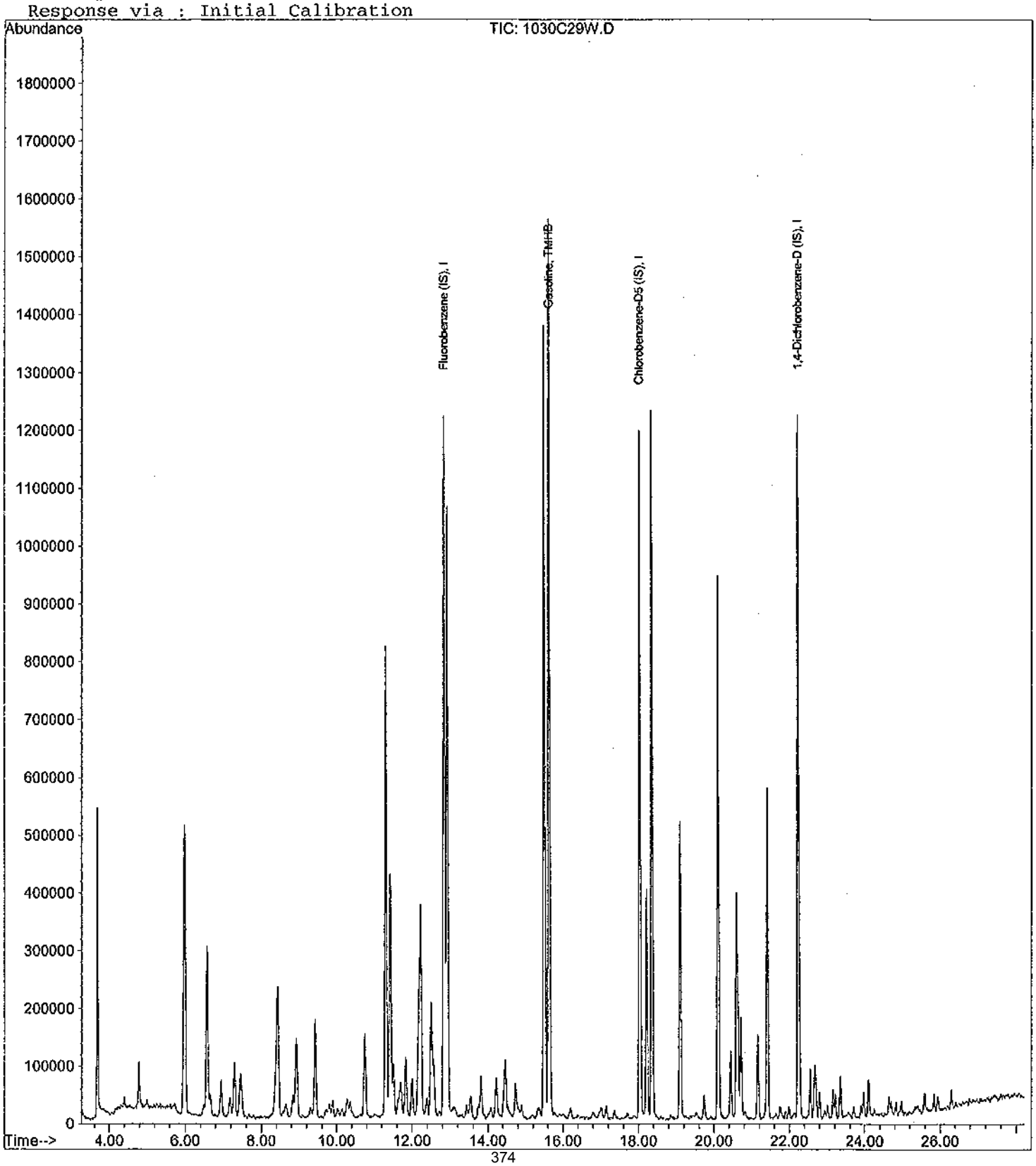
(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : M:\CHICO\DATA\C111030\1030C29W.D Vial: 1
Acq On : 31 Oct 11 9:31 Operator: STC
Sample : GAS 300ug/L (SS) Inst : Chico
Misc : Water 10mLw/ IS&S:10-30/10-26-11 Multiplr: 1.00

Quant Time: Nov 3 10:51 2011 Quant Results File: CGAS.RES

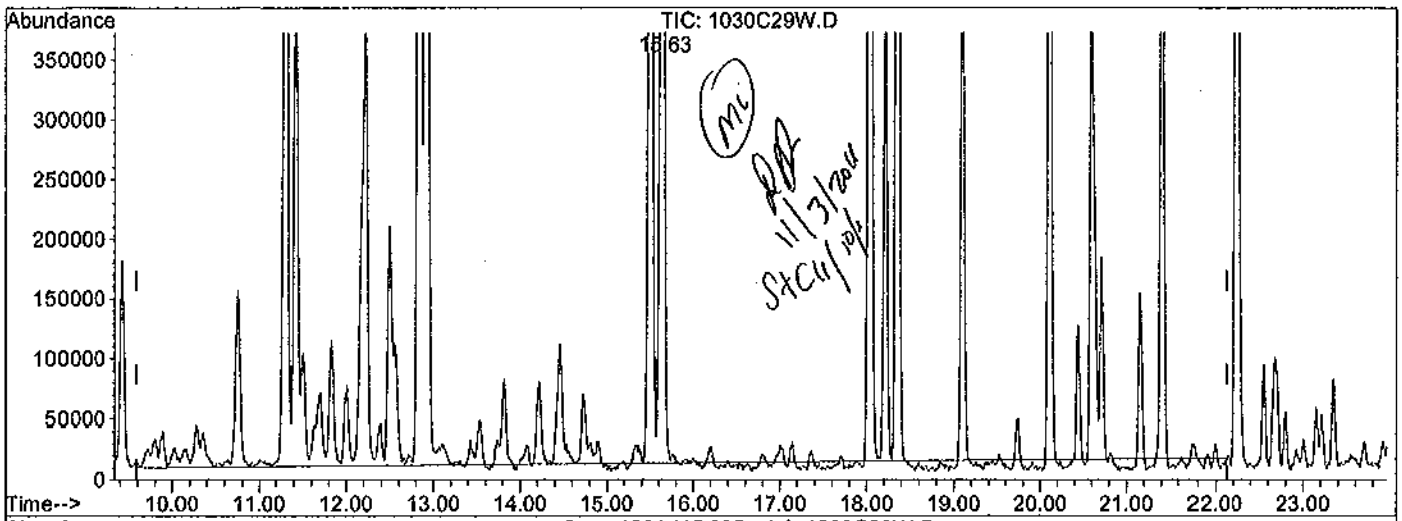
Method : M:\CHICO\DATA\C111030\CGAS.M (RTE Integrator)
Title : METHOD 8260B
Last Update : Thu Nov 03 10:47:02 2011
Response via : Initial Calibration

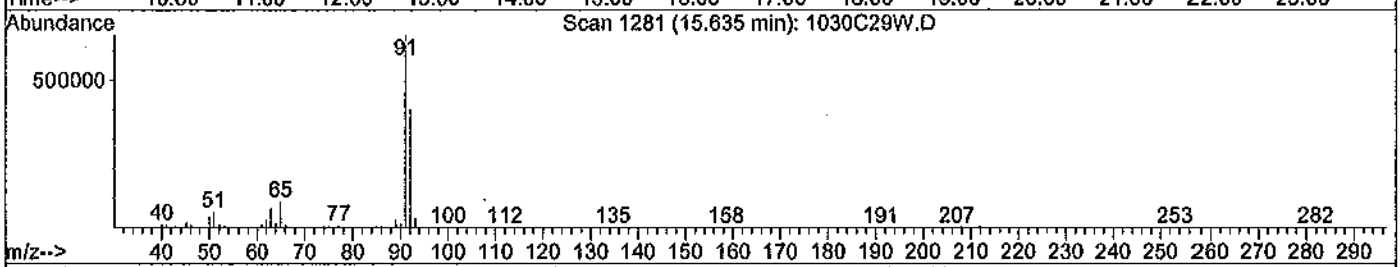


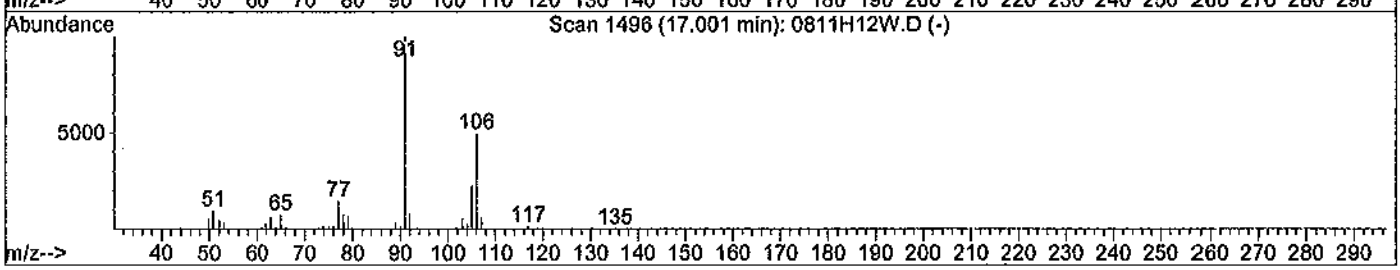
Quantitation Report

Data File : M:\CHICO\DATA\C111030\1030C29W.D Vial: 1
Acq On : 31 Oct 11 9:31 Operator: STC
Sample : GAS 300ug/L (SS) Inst : Chico
Misc : Water 10mLw/ IS&S:10-30/10-26-11 Multiplr: 1.00
Quant Time: Nov 3 10:47 2011 Quant Results File: temp.res

Method : M:\CHICO\DATA\C111030\CGAS.M (RTE Integrator)
Title : METHOD 8260B
Last Update : Thu Nov 03 10:47:02 2011
Response via : Multiple Level Calibration







TIC: 1030C29W.D

(2) Gasoline (TMHB)

15.64min 275.5469ppb m

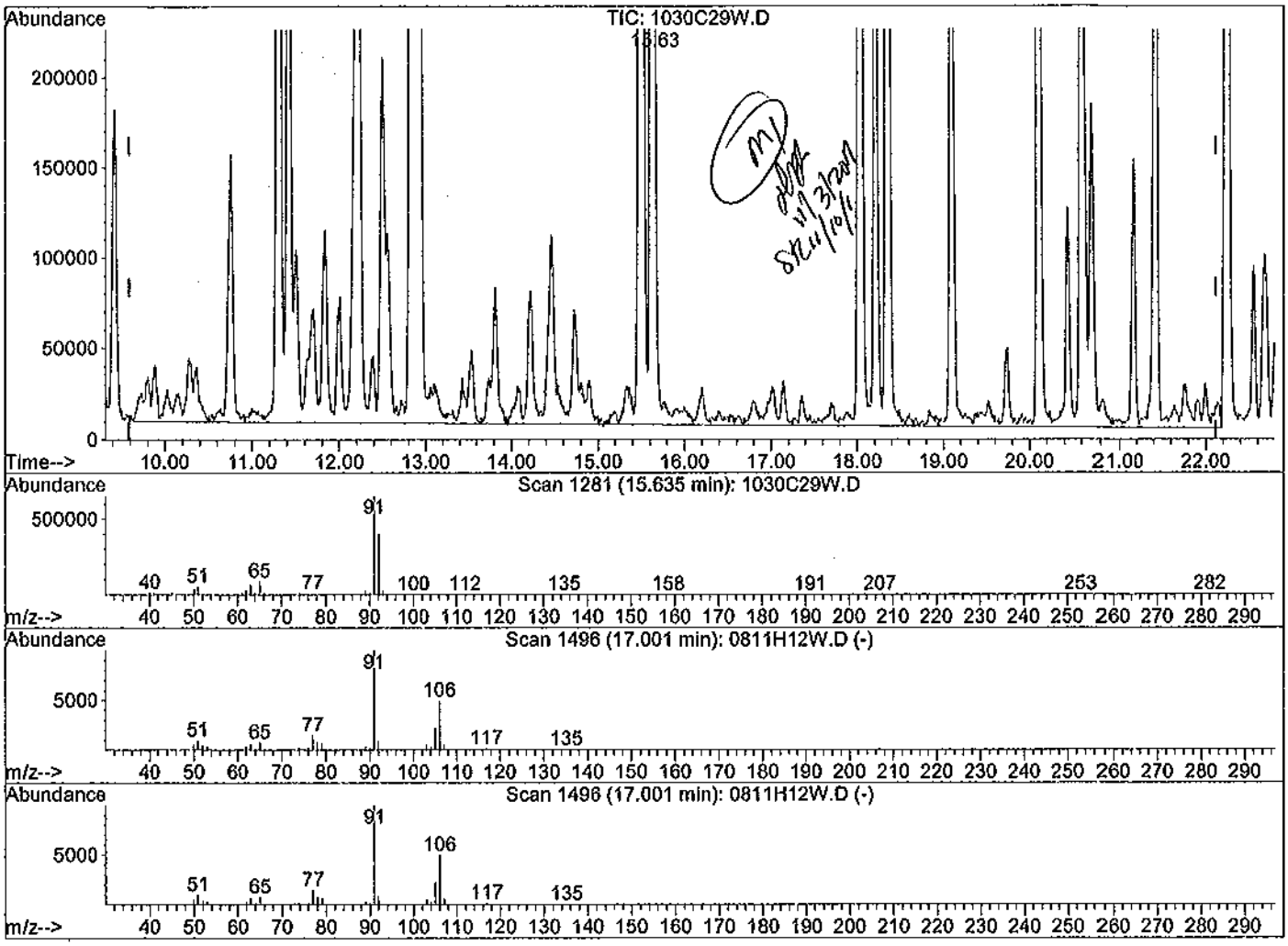
response 41492142

Ion	Exp%	Act%
TIC	100	100
0.00	0.00	0.26#
0.00	0.00	0.78#
0.00	0.00	0.00

Quantitation Report

Data File : M:\CHICO\DATA\C111030\1030C29W.D Vial: 1
Acq On : 31 Oct 11 9:31 Operator: STC
Sample : GAS 300ug/L (SS) Inst : Chico
Misc : Water 10mLw/ IS&S:10-30/10-26-11 Multiplr: 1.00
Quant Time: Nov 3 10:51 2011 Quant Results File: temp.res

Method : M:\CHICO\DATA\C111030\CGAS.M (RTE Integrator)
Title : METHOD 8260B
Last Update : Thu Nov 03 10:47:02 2011
Response via : Multiple Level Calibration



TIC: 1030C29W.D

(2) Gasoline (TMHB)

15.63min 332.6619ppb m

response 46900368

Ion	Exp%	Act%
TIC	100	100
0.00	0.00	0.23#
0.00	0.00	0.69#
0.00	0.00	0.00

VOLATILE ORGANIC ANALYSIS BY
EPA METHOD 8260B

Form 7

Continuing Calibration

Lab Name: APPL, Inc.
Case No:
Matrix:

SDG No:
Date Analyzed: 10/31/2011
Instrument: Chico
Initial Cal. Date: 10/30/2011
Data File: 1030C30W.D

		Compound	MEAN	CCRF	%D		%Drift
1	I	Fluorobenzene (IS)	ISTD			I	
2	TMHB	Gasoline	5.897	3.225	45	TMHBL	11
3	I	Chlorobenzene-D5 (IS)	ISTD			I	
4	I	1,4-Dichlorobenzene-D (IS)	ISTD			I	
5							
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36							
37							
38							
39							
40							
		Average			45.0		

Quantitation Report (Not Reviewed)

Data File : M:\CHICO\DATA\C111030\1030C30W.D Vial: 1
Acq On : 31 Oct 11 10:14 Operator: STC
Sample : GAS 300ug/L CCV-1WC Inst : Chico
Misc : Water 10mLw/ IS&S:10-30/10-26-11 Multiplr: 1.00

Quant Time: Nov 3 10:51 2011 Quant Results File: CGAS.RES

Quant Method : M:\CHICO\DATA\C111030\CGAS.M (RTE Integrator)
Title : METHOD 8260B
Last Update : Thu Nov 03 10:47:02 2011
Response via : Initial Calibration
DataAcq Meth : V8260

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Fluorobenzene (IS)	12.84	TIC	1207524	25.00000	ppb	0.00
3) Chlorobenzene-D5 (IS)	18.03	TIC	1180683	25.00000	ppb	0.00
4) 1,4-Dichlorobenzene-D (IS)	22.24	TIC	1196096	25.00000	ppb	0.00

System Monitoring Compounds

Target Compounds						Qvalue
2) Gasoline	15.64	TIC	46736698m	332.52710	ppb	100

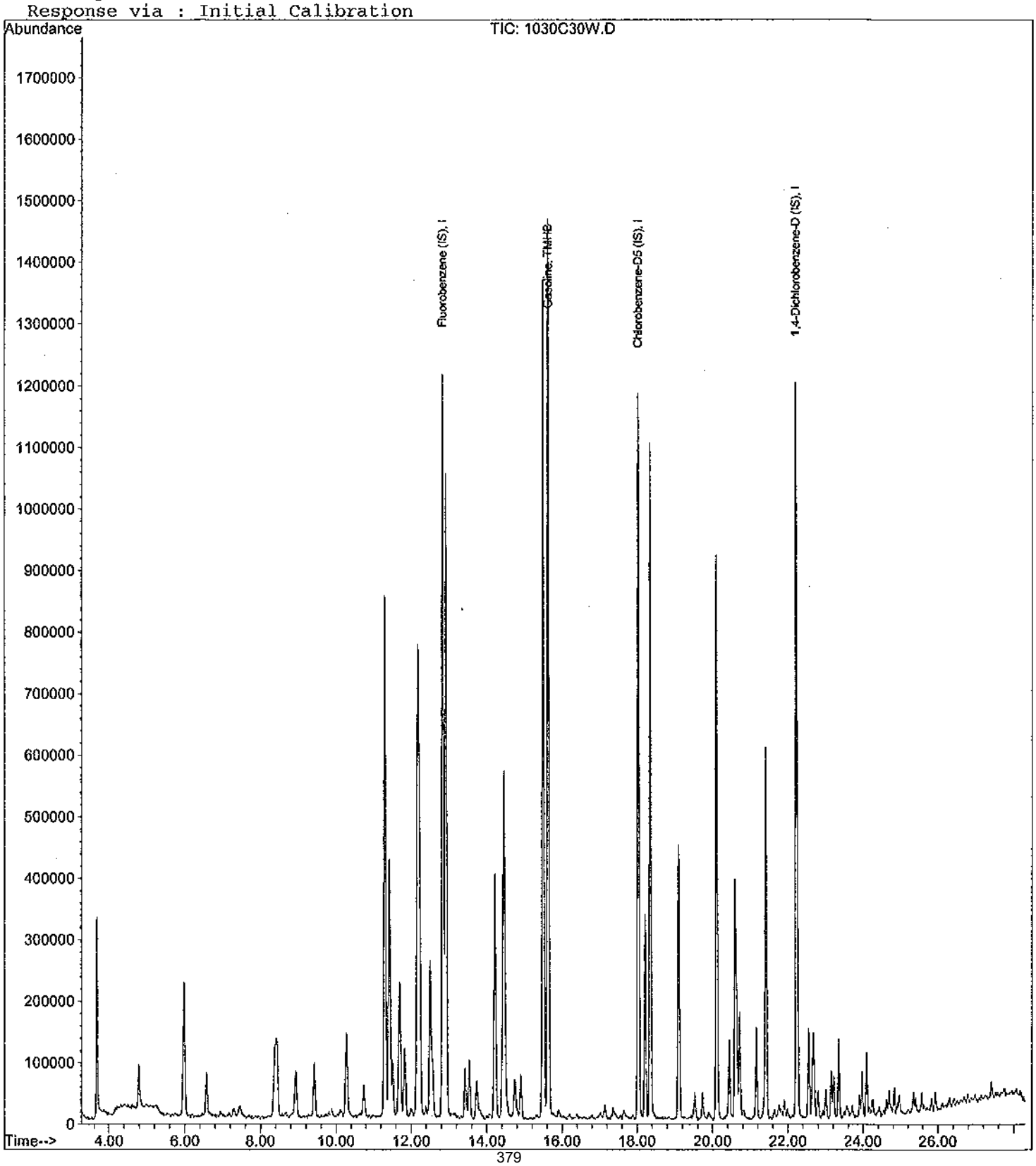
(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : M:\CHICO\DATA\C111030\1030C30W.D Vial: 1
Acq On : 31 Oct 11 10:14 Operator: STC
Sample : GAS 300ug/L CCV-1WC Inst : Chico
Misc : Water 10mLw/ IS&S:10-30/10-26-11 Multiplr: 1.00

Quant Time: Nov 3 10:51 2011 Quant Results File: CGAS.RES

Method : M:\CHICO\DATA\C111030\CGAS.M (RTE Integrator)
Title : METHOD 8260B
Last Update : Thu Nov 03 10:47:02 2011
Response via : Initial Calibration



EPA METHOD 8260B
Volatile Organic Compounds
Raw Data

APPL, INC.

Method Blank

EPA 8260B VOCs + Gas Water

Blank Name/QCG: **111030W-49334 - 161029**
Batch ID: #86RHB-111030AC

APPL Inc.
908 North Temperance Avenu
Clovis, CA 93611

Sample Type	Analyte	Result	LOQ	LOD	DL	Units	Extraction Date	Analysis Date
BLANK	1,1,1,2-TETRACHLOROETHANE	0.26 U	1.0	0.26	0.13	ug/L	10/31/11	10/31/11
BLANK	1,1,1-TRICHLOROETHANE	0.28 U	1.0	0.28	0.14	ug/L	10/31/11	10/31/11
BLANK	1,1,2,2-TETRACHLOROETHANE	0.20 U	1.0	0.20	0.10	ug/L	10/31/11	10/31/11
BLANK	1,1,2-TRICHLOROETHANE	0.40 U	1.0	0.40	0.20	ug/L	10/31/11	10/31/11
BLANK	1,1-DICHLOROETHANE	0.38 U	1.0	0.38	0.19	ug/L	10/31/11	10/31/11
BLANK	1,1-DICHLOROETHENE	0.60 U	1.0	0.60	0.30	ug/L	10/31/11	10/31/11
BLANK	1,2,3-TRICHLOROPROPANE	0.78 U	2.0	0.78	0.39	ug/L	10/31/11	10/31/11
BLANK	1,2,4-TRICHLOROBENZENE	0.42 U	1.0	0.42	0.21	ug/L	10/31/11	10/31/11
BLANK	1,2-DIBROMO-3-CHLOROPROPA	1.52 U	2.0	1.52	0.76	ug/L	10/31/11	10/31/11
BLANK	1,2-DIBROMOETHANE	0.40 U	1.0	0.40	0.20	ug/L	10/31/11	10/31/11
BLANK	1,2-DICHLOROBENZENE	0.34 U	1.0	0.34	0.17	ug/L	10/31/11	10/31/11
BLANK	1,2-DICHLOROETHANE	0.28 U	1.0	0.28	0.14	ug/L	10/31/11	10/31/11
BLANK	1,2-DICHLOROPROPANE	0.34 U	1.0	0.34	0.17	ug/L	10/31/11	10/31/11
BLANK	1,3-DICHLOROBENZENE	0.22 U	1.0	0.22	0.11	ug/L	10/31/11	10/31/11
BLANK	1,3-DICHLOROPROPENE, TOTAL	0.36 U	1.0	0.36	0.18	ug/L	10/31/11	10/31/11
BLANK	1,4-DICHLOROBENZENE	0.38 U	1.0	0.38	0.19	ug/L	10/31/11	10/31/11
BLANK	2-BUTANONE	1.20 U	10.0	1.20	0.60	ug/L	10/31/11	10/31/11
BLANK	4-METHYL-2-PENTANONE	3.80 U	10.0	3.80	1.90	ug/L	10/31/11	10/31/11
BLANK	ACETONE	1.90 U	10.0	1.90	0.95	ug/L	10/31/11	10/31/11
BLANK	BENZENE	0.32 U	1.0	0.32	0.16	ug/L	10/31/11	10/31/11
BLANK	BROMODICHLOROMETHANE	0.28 U	1.0	0.28	0.14	ug/L	10/31/11	10/31/11
BLANK	BROMOFORM	0.28 U	1.0	0.28	0.14	ug/L	10/31/11	10/31/11
BLANK	BROMOMETHANE	0.48 U	2.0	0.48	0.24	ug/L	10/31/11	10/31/11
BLANK	CARBON TETRACHLORIDE	0.20 U	1.0	0.20	0.10	ug/L	10/31/11	10/31/11
BLANK	CHLOROBENZENE	0.42 U	1.0	0.42	0.21	ug/L	10/31/11	10/31/11
BLANK	CHLORODIBROMOMETHANE	0.38 U	1.0	0.38	0.19	ug/L	10/31/11	10/31/11
BLANK	CHLOROETHANE	0.42 U	1.0	0.42	0.21	ug/L	10/31/11	10/31/11
BLANK	CHLOROFORM	0.14 U	1.0	0.14	0.07	ug/L	10/31/11	10/31/11
BLANK	CHLOROMETHANE	0.84 U	1.0	0.84	0.42	ug/L	10/31/11	10/31/11
BLANK	CIS-1,2-DICHLOROETHENE	0.32 U	1.0	0.32	0.16	ug/L	10/31/11	10/31/11
BLANK	ETHYLBENZENE	0.46 U	1.0	0.46	0.23	ug/L	10/31/11	10/31/11
BLANK	GASOLINE	12.12 U	20.0	12.12	6.06	ug/L	10/31/11	10/31/11
BLANK	HEXACHLOROBUTADIENE	0.38 U	1.0	0.38	0.19	ug/L	10/31/11	10/31/11
BLANK	METHYL TERT-BUTYL ETHER	0.52 U	1.0	0.52	0.26	ug/L	10/31/11	10/31/11

Quant Method:CALLW.M
Run #:1030C34
Instrument:Chico
Sequence:C111030
Initials:ARS

GC SC-Blank-REG MDLs
Printed: 12/06/11 3:06:21 PM

Method Blank
EPA 8260B VOCs + Gas Water

Blank Name/QCG: **111030W-49334 - 161029**
Batch ID: #86RHB-111030AC

APPL Inc.
908 North Temperance Avenu
Clovis, CA 93611

Sample Type	Analyte	Result	LOQ	LOD	DL	Units	Extraction Date	Analysis Date
BLANK	METHYLENE CHLORIDE	0.70 U	5.0	0.70	0.35	ug/L	10/31/11	10/31/11
BLANK	STYRENE	0.50 U	1.0	0.50	0.25	ug/L	10/31/11	10/31/11
BLANK	TETRACHLOROETHENE	0.48 U	1.0	0.48	0.24	ug/L	10/31/11	10/31/11
BLANK	TOLUENE	0.34 U	1.0	0.34	0.17	ug/L	10/31/11	10/31/11
BLANK	TRANS-1,2-DICHLOROETHENE	0.38 U	1.0	0.38	0.19	ug/L	10/31/11	10/31/11
BLANK	TRICHLOROETHENE	0.32 U	1.0	0.32	0.16	ug/L	10/31/11	10/31/11
BLANK	VINYL CHLORIDE	0.46 U	1.0	0.46	0.23	ug/L	10/31/11	10/31/11
BLANK	XYLENES (TOTAL)	0.38 U	1.0	0.38	0.19	ug/L	10/31/11	10/31/11
BLANK	SURROGATE: 1,2-DICHLOROET	109	70-120			%	10/31/11	10/31/11
BLANK	SURROGATE: 4-BROMOFLUORO	102	75-120			%	10/31/11	10/31/11
BLANK	SURROGATE: DIBROMOFLUOR	105	85-115			%	10/31/11	10/31/11
BLANK	SURROGATE: TOLUENE-D8 (S)	102	85-120			%	10/31/11	10/31/11

Quant Method:CALLW.M
Run #:1030C34
Instrument:Chico
Sequence:C111030
Initials:ARS

GC SC-Blank-REG MDLs
Printed: 12/06/11 3:06:21 PM

Quantitation Report (QT Reviewed)

Data File : M:\CHICO\DATA\C111030\1030C34W.D Vial: 1
Acq On : 31 Oct 11 13:02 Operator: STC
Sample : 111030A BLk-1WC Inst : Chico
Misc : Water 10mLw/ IS&S:10-30/10-26-11 Multiplr: 1.00

Quant Time: Nov 3 11:50 2011 Quant Results File: CALLW.RES

Quant Method : M:\CHICO\DATA\C111030\CALLW.M (RTE Integrator)
Title : METHOD 8260
Last Update : Wed Nov 02 14:33:25 2011
Response via : Initial Calibration
DataAcq Meth : V8260

Internal Standards	R.T.	QIon	Response	Conc Units	Dev(Min)
1) Fluorobenzene (IS)	12.85	96	564158	25.00000 ppb	0.02
55) Chlorobenzene-D5 (IS)	18.05	117	392640	25.00000 ppb	0.00
71) 1,4-Dichlorobenzene-D (IS)	22.26	152	207424	25.00000 ppb	0.02
System Monitoring Compounds					
33) Dibromofluoromethane(S)	11.43	111	396364	26.37439 ppb	0.00
Spiked Amount 25.097			Recovery = 105.087%		
38) 1,2-DCA-D4(S)	12.23	65	352223	26.32879 ppb	0.00
Spiked Amount 24.225			Recovery = 108.684%		
56) Toluene-D8(S)	15.51	98	1448391	26.21699 ppb	0.02
Spiked Amount 25.808			Recovery = 101.584%		
64) 4-Bromofluorobenzene(S)	20.12	95	516404	26.08735 ppb	0.00
Spiked Amount 25.459			Recovery = 102.465%		

Target Compounds Qvalue

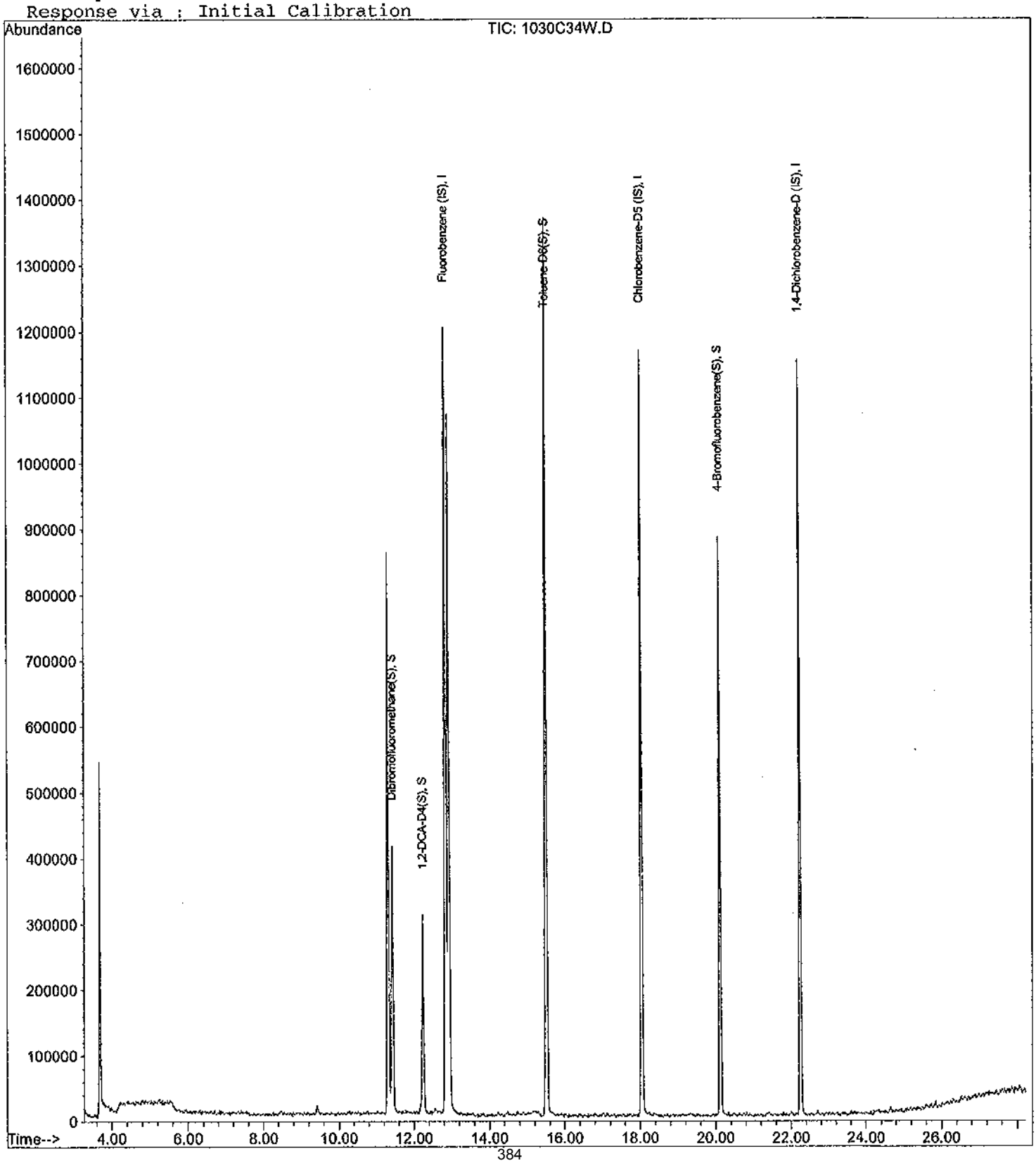
(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : M:\CHICO\DATA\C111030\1030C34W.D Vial: 1
Acq On : 31 Oct 11 13:02 Operator: STC
Sample : 111030A BLk-1WC Inst : Chico
Misc : Water 10mLw/ IS&S:10-30/10-26-11 Multiplr: 1.00

Quant Time: Nov 3 11:50 2011 Quant Results File: CALLW.RES

Method : M:\CHICO\DATA\C111030\CALLW.M (RTE Integrator)
Title : METHOD 8260
Last Update : Thu Nov 03 10:27:07 2011
Response via : Initial Calibration



Quantitation Report (QT Reviewed)

Data File : M:\CHICO\DATA\C111030\1030C34W.D Vial: 1
Acq On : 31 Oct 11 13:02 Operator: STC
Sample : 111030A BLk-1WC Inst : Chico
Misc : Water 10mLw/ IS&S:10-30/10-26-11 Multiplr: 1.00

Quant Time: Nov 10 10:22 2011 Quant Results File: CGAS.RES

Quant Method : M:\CHICO\DATA\C111030\CGAS.M (RTE Integrator)
Title : METHOD 8260B
Last Update : Thu Nov 03 10:47:02 2011
Response via : Initial Calibration
DataAcq Meth : V8260

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Fluorobenzene (IS)	12.84	TIC	1192714	25.00000	ppb	0.00
3) Chlorobenzene-D5 (IS)	18.05	TIC	1166397	25.00000	ppb	0.01
4) 1,4-Dichlorobenzene-D (IS)	22.25	TIC	1149644	25.00000	ppb	0.00

System Monitoring Compounds

Target Compounds Qvalue

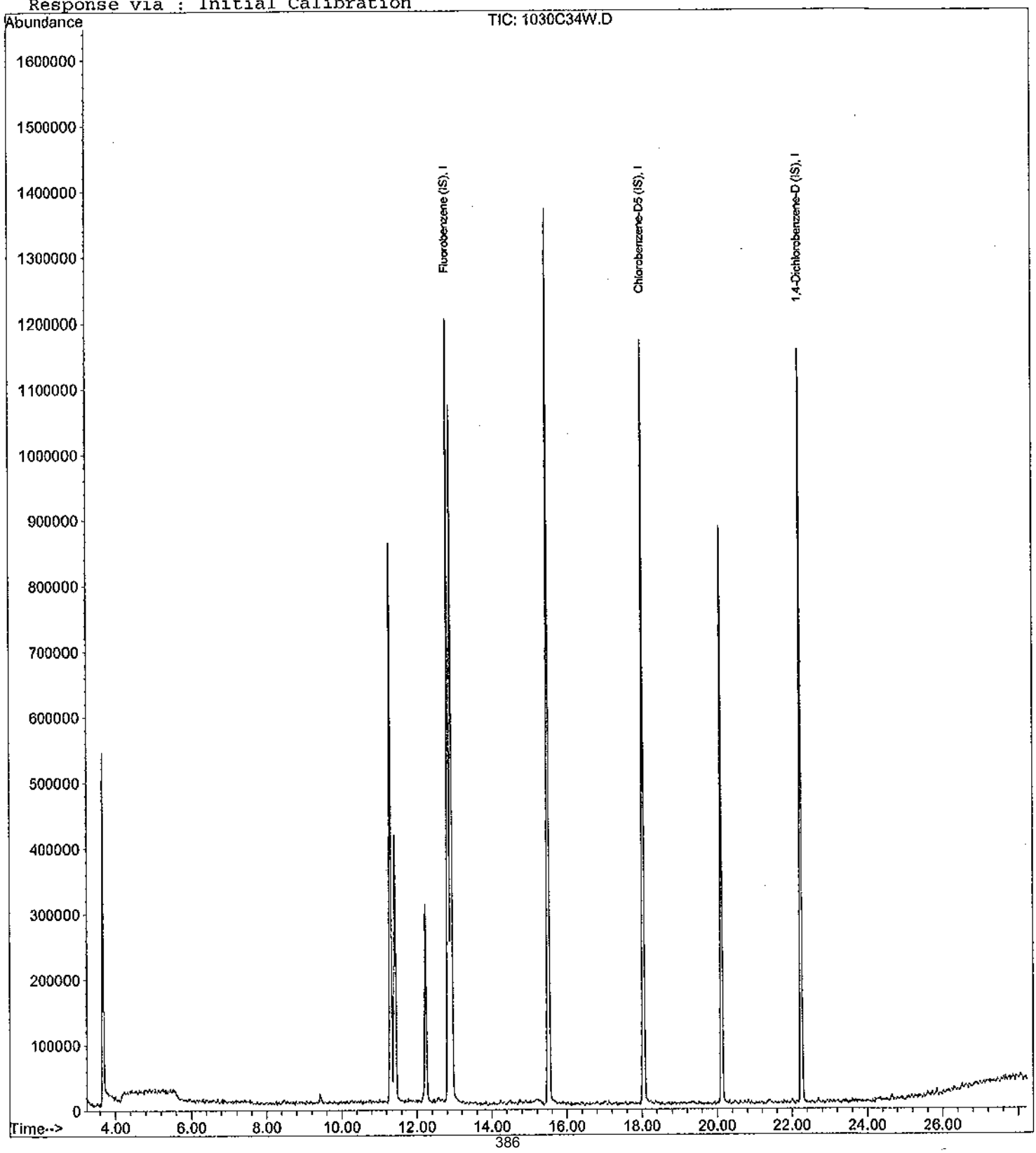
Quantitation Report

Data File : M:\CHICO\DATA\C111030\1030C34W.D
Acq On : 31 Oct 11 13:02
Sample : 111030A BLk-1WC
Misc : Water 10mLw/ IS&S:10-30/10-26-11

Vial: 1
Operator: STC
Inst : Chico
Multiplr: 1.00

Quant Time: Nov 10 10:22 2011 Quant Results File: CGAS.RES

Method : M:\CHICO\DATA\C111030\CGAS.M (RTE Integrator)
Title : METHOD 8260B
Last Update : Thu Nov 03 10:47:02 2011
Response via : Initial Calibration



Laboratory Control Spike Recovery
EPA 8260B VOCs + Gas Water

APPL ID: **111031W-49334 LCS - 161029**
Batch ID: #86RHB-111030AC

APPL Inc.
908 North Temperance Avenue
Clovis, CA 93611

Compound Name	Spike Level ug/L	SPK Result ug/L	SPK % Recovery	Recovery Limits
1,1,1,2-TETRACHLOROETHANE	10.00	11.0	110	80-130
1,1,1-TRICHLOROETHANE	10.00	9.55	95.5	65-130
1,1,2,2-TETRACHLOROETHANE	10.00	10.3	103	65-130
1,1,2-TRICHLOROETHANE	10.00	9.91	99.1	75-125
1,1-DICHLOROETHANE	10.00	9.98	99.8	70-135
1,1-DICHLOROETHENE	10.00	8.83	88.3	70-130
1,2,3-TRICHLOROPROPANE	10.00	11.8	118	75-125
1,2,4-TRICHLOROBENZENE	10.00	10.2	102	65-135
1,2-DIBROMO-3-CHLOROPROPANE	10.00	10.2	102	50-130
1,2-DIBROMOETHANE	10.00	10.7	107	70-130
1,2-DICHLOROBENZENE	10.00	10.1	101	70-120
1,2-DICHLOROETHANE	10.00	9.61	96.1	70-130
1,2-DICHLOROPROPANE	10.00	10.1	101	75-125
1,3-DICHLOROBENZENE	10.00	9.67	96.7	75-125
1,3-DICHLOROPROPENE, TOTAL	20.0	19.7	98.5	70-130
1,4-DICHLOROBENZENE	10.00	9.94	99.4	75-125
2-BUTANONE	10.00	10.5	105	30-150
4-METHYL-2-PENTANONE	10.00	9.78	97.8	60-135
ACETONE	10.00	13.0	130	40-140
BENZENE	10.00	9.48	94.8	80-120
BROMODICHLOROMETHANE	10.00	10.0	100	75-120
BROMOFORM	10.00	9.32	93.2	70-130
BROMOMETHANE	10.00	8.95	89.5	30-145
CARBON TETRACHLORIDE	10.00	9.77	97.7	65-140
CHLOROBENZENE	10.00	10.3	103	80-120
CHLORODIBROMOMETHANE	10.00	11.0	110	60-135

Comments:

Primary	SPK
Quant Method :	CALLW.M
Extraction Date :	10/31/11
Analysis Date :	10/31/11
Instrument :	Chico
Run :	1030C28
Initials :	ARS

Printed: 12/06/11 3:06:22 PM

APPL Standard LCS

Laboratory Control Spike Recovery
EPA 8260B VOCs + Gas Water

APPL ID: **111031W-49334 LCS - 161029**
Batch ID: #86RHB-111030AC

APPL Inc.
908 North Temperance Avenue
Clovis, CA 93611

Compound Name	Spike Level ug/L	SPK Result ug/L	SPK % Recovery	Recovery Limits
CHLOROETHANE	10.00	8.89	88.9	60-135
CHLOROFORM	10.00	9.79	97.9	65-135
CHLOROMETHANE	10.00	9.42	94.2	40-125
CIS-1,2-DICHLOROETHENE	10.00	9.18	91.8	70-125
ETHYLBENZENE	10.00	10.1	101	75-125
GASOLINE	300	370	123	75-125
HEXACHLOROBUTADIENE	10.00	10.5	105	50-140
METHYL TERT-BUTYL ETHER	10.00	9.69	96.9	65-125
METHYLENE CHLORIDE	10.00	9.05	90.5	55-140
STYRENE	10.00	10.6	106	65-135
TETRACHLOROETHENE	10.00	10.1	101	45-150
TOLUENE	10.00	9.40	94.0	75-120
TRANS-1,2-DICHLOROETHENE	10.00	9.05	90.5	60-140
TRICHLOROETHENE	10.00	9.64	96.4	70-125
VINYL CHLORIDE	10.00	10.2	102	50-145
XYLENES (TOTAL)	30.0	30.3	101	80-120
SURROGATE: 1,2-DICHLOROETHANE-D	24.2	24.4	101	70-120
SURROGATE: 4-BROMOFLUOROBENZE	25.5	27.7	109	75-120
SURROGATE: DIBROMOFLUOROMETH	25.1	25.2	100	85-115
SURROGATE: TOLUENE-D8 (S)	25.8	26.9	104	85-120

Comments:

Primary	SPK
Quant Method :	CALLW.M
Extraction Date :	10/31/11
Analysis Date :	10/31/11
Instrument :	Chlco
Run :	1030C28
Initials :	ARS

Printed: 12/06/11 3:06:22 PM

APPL Standard LCS

Quantitation Report (Not Reviewed)

Data File : M:\CHICO\DATA\C111030\1030C28W.D Vial: 1
Acq On : 31 Oct 11 8:48 Operator: STC
Sample : 111030A LCS-1WC (SS) Inst : Chico
Misc : Water 10mLw/ IS&S:10-30/10-26-11 Multiplr: 1.00

Quant Time: Dec 2 11:35 2011 Quant Results File: CALLW.RES

Quant Method : M:\CHICO\DATA\C111030\CALLW.M (RTE Integrator)
Title : METHOD 8260
Last Update : Fri Dec 02 11:32:50 2011
Response via : Initial Calibration
DataAcq Meth : V8260

Internal Standards	R.T.	QIon	Response	Conc	Units		Dev(Min)
1) Fluorobenzene (IS)	12.84	96	600576	25.00000	ppb		0.00
55) Chlorobenzene-D5 (IS)	18.04	117	389760	25.00000	ppb		0.00
71) 1,4-Dichlorobenzene-D (IS)	22.24	152	212800	25.00000	ppb		0.00
System Monitoring Compounds							
33) Dibromofluoromethane(S)	11.42	111	402364	25.15012	ppb		0.00
Spiked Amount 25.097			Recovery = 100.210%				
38) 1,2-DCA-D4(S)	12.23	65	347346	24.38980	ppb		0.00
Spiked Amount 24.225			Recovery = 100.680%				
56) Toluene-D8(S)	15.50	98	1474138	26.88019	ppb		0.00
Spiked Amount 25.808			Recovery = 104.153%				
64) 4-Bromofluorobenzene(S)	20.11	95	543410	27.65447	ppb		0.00
Spiked Amount 25.459			Recovery = 108.620%				
Target Compounds							Qvalue
2) Dichlorodifluoromethane	4.07	85	222193	10.04166	ppb		100
3) Freon 114	4.33	85	146402	10.56428	ppb		93
4) Chloromethane	4.55	50	258385	9.42321	ppb		99
5) Vinyl chloride	4.82	62	186738	10.18146	ppb		96
7) Bromomethane	5.72	94	119189	8.95228	ppb		94
8) Chloroethane	5.91	64	134747	8.89428	ppb		97
9) Dichlorofluoromethane	6.01	67	386134	9.21828	ppb		97
10) Trichlorofluoromethane	6.52	101	241622	9.72027	ppb		99
11) Acetonitrile	7.64	41	78885	119.90087	ug/l		100
12) Acrolein	7.16	56	34469	114.48914	ppb		96
13) Acetone	7.27	43	22365	12.99757	ppb	#	84
14) Freon-113	7.46	101	138327	9.44782	ppb		97
15) 1,1-DCE	7.67	96	151407	8.83040	ppb		96
16) t-Butanol	7.76	59	10529	129.32077	ppb		93
17) Methyl Acetate	8.18	43	48755	9.36519	ppb		96
18) Iodomethane	8.16	142	102169	10.69989	ppb		90
19) Acrylonitrile	8.56	53	17916	9.49044	ppb		79
20) Methylene chloride	8.47	84	147951	9.04673	ppb		99
21) Carbon disulfide	8.56	76	154944	9.30051	ppb		100
22) Methyl t-butyl ether (MtBE	8.89	73	251165	9.69151	ppb		96
23) Trans-1,2-DCE	9.10	96	180083	9.05360	ppb		88
24) Diisopropyl Ether	9.75	45	553904	9.66784	ppb		94
25) 1,1-DCA	9.79	63	339012	9.97748	ppb		99
26) Vinyl Acetate	9.42	43	104836	9.85383	ppb	#	83
27) Ethyl tert Butyl Ether	10.45	59	395408	10.10888	ppb		99
28) MEK (2-Butanone)	10.44	43	71405	10.48433	ppb		99
29) Cis-1,2-DCE	10.82	96	187663	9.18084	ppb		97
30) 2,2-Dichloropropane	10.82	77	208247	8.55600	ppb		97
31) Chloroform	11.10	83	320091	9.79258	ppb		99
32) Bromochloromethane	11.32	128	58472	10.27501	ppb		98
34) 1,1,1-TCA	11.84	97	283983	9.55329	ppb		96
35) Cyclohexane	12.00	56	268948	9.71733	ppb		94
36) 1,1-Dichloropropene	12.10	75	240188	9.42981	ppb		99
37) 2,2,4-Trimethylpentane	12.18	57	414455	9.47423	ppb		98
39) Carbon Tetrachloride	12.30	117	199898	9.76577	ppb		98
40) Tert Amyl Methyl Ether	12.34	73	292021	9.98459	ppb		99

(#) = qualifier out of range (m) = manual integration

Quantitation Report (Not Reviewed)

Data File : M:\CHICO\DATA\C111030\1030C28W.D Vial: 1
Acq On : 31 Oct 11 8:48 Operator: STC
Sample : 111030A LCS-1WC (SS) Inst : Chico
Misc : Water 10mLw/ IS&S:10-30/10-26-11 Multiplr: 1.00

Quant Time: Dec 2 11:35 2011 Quant Results File: CALLW.RES

Quant Method : M:\CHICO\DATA\C111030\CALLW.M (RTE Integrator)
Title : METHOD 8260
Last Update : Fri Dec 02 11:32:50 2011
Response via : Initial Calibration
DataAcq Meth : V8260

	Compound	R.T.	QIon	Response	Conc	Unit		Qvalue
41)	1,2-DCA	12.38	62	161160	9.60516	ppb		100
42)	Benzene	12.50	78	693647	9.47863	ppb		96
43)	TCE	13.53	95	195399	9.64105	ppb		91
44)	2-Pentanone	13.20	43	531259	125.31057	ppb		97
45)	1,2-Dichloropropane	13.76	63	167281	10.05687	ppb	#	94
46)	Bromodichloromethane	14.11	83	190021	9.99949	ppb	#	91
47)	Methyl Cyclohexane	13.82	83	224538	9.48040	ppb		99
48)	Dibromomethane	14.16	93	67527	10.15102	ppb		95
49)	2-Chloroethyl vinyl ether	14.57	63	42382	10.02259	ppb		95
50)	1-Bromo-2-chloroethane	14.88	63	147231	10.37368	ppb	#	79
51)	Cis-1,3-Dichloropropene	15.00	75	181879	10.03716	ppb		100
52)	Toluene	15.63	91	678338	9.39804	ppb		99
53)	Trans-1,3-Dichloropropene	15.80	75	126600	9.70449	ppb		98
54)	1,1,2-TCA	16.08	83	69681	9.91034	ppb		93
57)	1,2-EDB	17.33	107	79107	10.66417	ppb		94
58)	Tetrachloroethene	16.78	164	202010	10.08114	ppb		94
59)	1-Chlorohexane	17.70	91	232734	10.08580	ppb		97
60)	1,1,1,2-Tetrachloroethane	18.16	131	138607	11.04877	ppb		99
61)	m&p-Xylene	18.35	106	589147	19.90172	ppb		97
62)	o-Xylene	19.11	106	295217	10.36928	ppb		98
63)	Styrene	19.13	104	457607	10.64883	ppb		93
65)	2-Hexanone	16.11	43	38770	10.87089	ppb		95
66)	1,3-Dichloropropane	16.49	76	147530	10.08561	ppb		98
67)	Dibromochloromethane	16.97	129	105397	11.03714	ppb		82
68)	Chlorobenzene	18.10	112	437982	10.34243	ppb		97
69)	Ethylbenzene	18.22	91	794180	10.07104	ppb		100
70)	Bromoform	19.65	173	45131	9.31734	ppb		91
72)	MIBK (methyl isobutyl keto	14.68	43	59150	9.77763	ppb		87
73)	Isopropylbenzene	19.73	105	756513	9.80877	ppb		98
74)	1,1,2,2-Tetrachloroethane	19.90	83	66287	10.26718	ppb	#	74
75)	1,2,3-Trichloropropane	20.16	110	8565	11.81260	ppb		82
76)	t-1,4-Dichloro-2-Butene	20.23	53	14963	10.22116	ppb	#	92
77)	Bromobenzene	20.48	156	179052	10.06445	ppb		89
78)	n-Propylbenzene	20.44	91	900774	9.79012	ppb		100
79)	4-Ethyltoluene	20.63	105	593563	9.32229	ppb		97
80)	2-Chlorotoluene	20.74	91	598129	9.81561	ppb		98
81)	1,3,5-Trimethylbenzene	20.72	105	617840	9.86323	ppb		99
82)	4-Chlorotoluene	20.82	91	502123	9.56935	ppb		98
83)	Tert-Butylbenzene	21.36	119	672218	9.91209	ppb		97
84)	1,2,4-Trimethylbenzene	21.42	105	604092	9.23368	ppb		96
85)	Sec-Butylbenzene	21.76	105	823845	10.12964	ppb		96
86)	p-Isopropyltoluene	21.99	119	683604	9.81315	ppb		98
87)	Benzyl Chloride	22.42	91	81362	8.79846	ppb		94
88)	1,3-DCB	22.12	146	351790	9.66926	ppb		95
89)	1,4-DCB	22.30	146	335795	9.94420	ppb		96
90)	Hexachloroethane	23.59	117	92345	8.87470	ppb		87
91)	n-Butylbenzene	22.69	91	582962	9.59448	ppb		98
92)	1,2-DCB	22.93	146	292666	10.11316	ppb		97
93)	1,2-Dibromo-3-chloropropan	24.14	155	10559	10.20135	ppb		92
94)	1,2,4-Trichlorobenzene	25.59	180	213173	10.16551	ppb		98
95)	Hexachlorobutadiene	25.84	223	40176	10.54513	ppb		97

(#) = qualifier out of range (m) = manual integration

Quantitation Report (Not Reviewed)

Data File : M:\CHICO\DATA\C111030\1030C28W.D Vial: 1
Acq On : 31 Oct 11 8:48 Operator: STC
Sample : 111030A LCS-1WC (SS) Inst : Chico
Misc : Water 10mLw/ IS&S:10-30/10-26-11 Multiplr: 1.00

Quant Time: Dec 2 11:35 2011 Quant Results File: CALLW.RES

Quant Method : M:\CHICO\DATA\C111030\CALLW.M (RTE Integrator)
Title : METHOD 8260
Last Update : Fri Dec 02 11:32:50 2011
Response via : Initial Calibration
DataAcq Meth : V8260

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
96) Naphthalene	25.94	128	272964	10.54986	ppb	99
97) 1,2,3-Trichlorobenzene	26.29	180	172357	10.86589	ppb	99

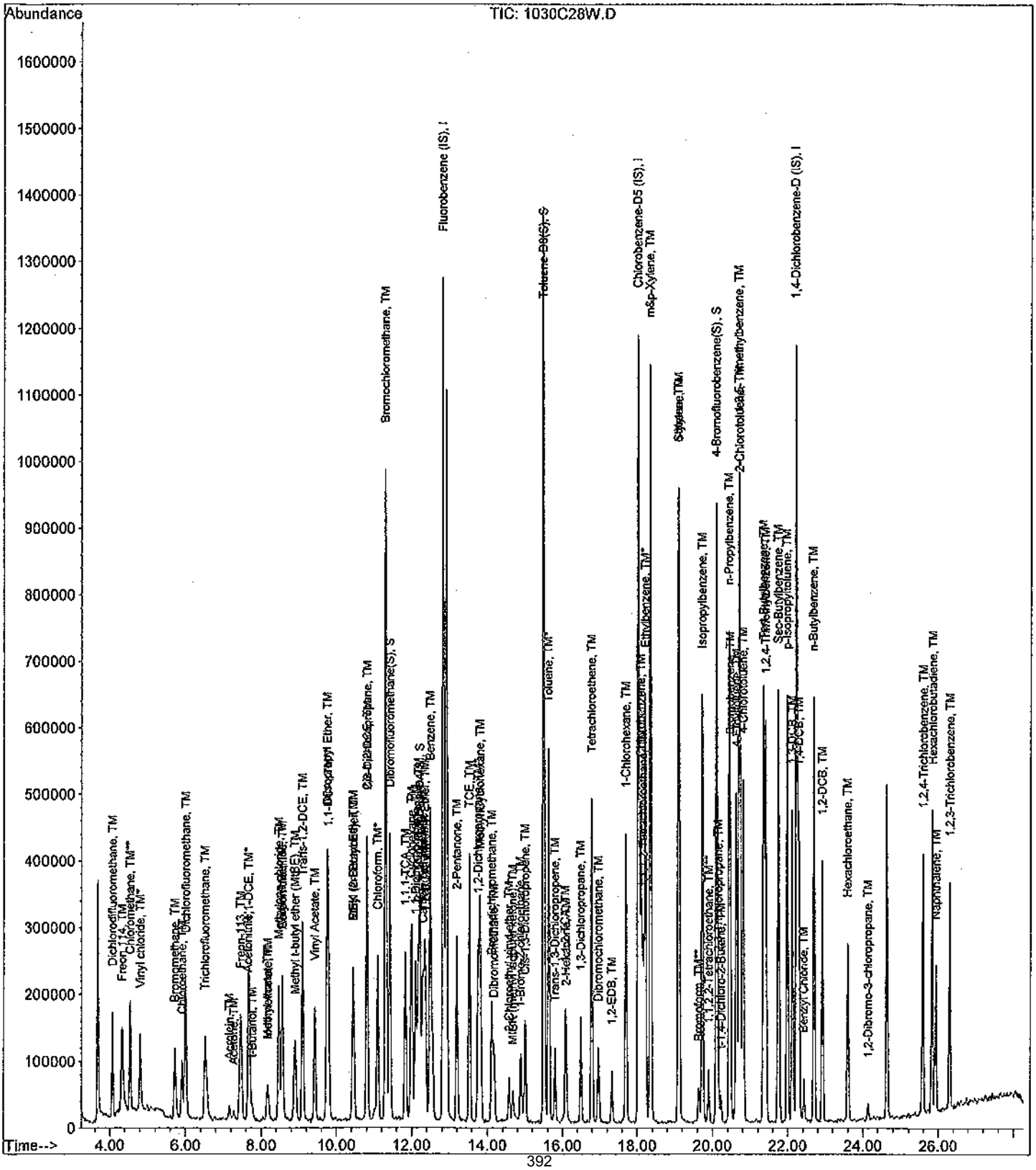
(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : M:\CHICO\DATA\C111030\1030C28W.D Vial: 1
Acq On : 31 Oct 11 8:48 Operator: STC
Sample : 111030A LCS-1WC (SS) Inst : Chico
Misc : Water 10mLw/ IS&S:10-30/10-26-11 Multiplr: 1.00

Quant Time: Dec 2 11:35 2011 Quant Results File: CALLW.RES

Method : M:\CHICO\DATA\C111030\CALLW.M (RTE Integrator)
Title : METHOD 8260
Last Update : Fri Dec 02 11:32:50 2011
Response via : Initial Calibration



Quantitation Report (QT Reviewed)

Data File : M:\CHICO\DATA\C111030\1030C31W.D Vial: 1
Acq On : 31 Oct 11 10:57 Operator: STC
Sample : GAS 300ug/L LCS-1WC Inst : Chico
Misc : Water 10mLw/ IS&S:10-30/10-26-11 Multiplr: 1.00

Quant Time: Nov 3 10:53 2011 Quant Results File: CGAS.RES

Quant Method : M:\CHICO\DATA\C111030\CGAS.M (RTE Integrator)
Title : METHOD 8260B
Last Update : Thu Nov 03 10:47:02 2011
Response via : Initial Calibration
DataAcq Meth : V8260

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Fluorobenzene (IS)	12.84	TIC	1238145	25.00000	ppb	0.00
3) Chlorobenzene-D5 (IS)	18.04	TIC	1217687	25.00000	ppb	0.00
4) 1,4-Dichlorobenzene-D (IS)	22.25	TIC	1183942	25.00000	ppb	0.00

System Monitoring Compounds

Target Compounds						Qvalue
2) Gasoline	15.64	TIC	51531735m	369.82724	ppb	100

(#) = qualifier out of range (m) = manual integration

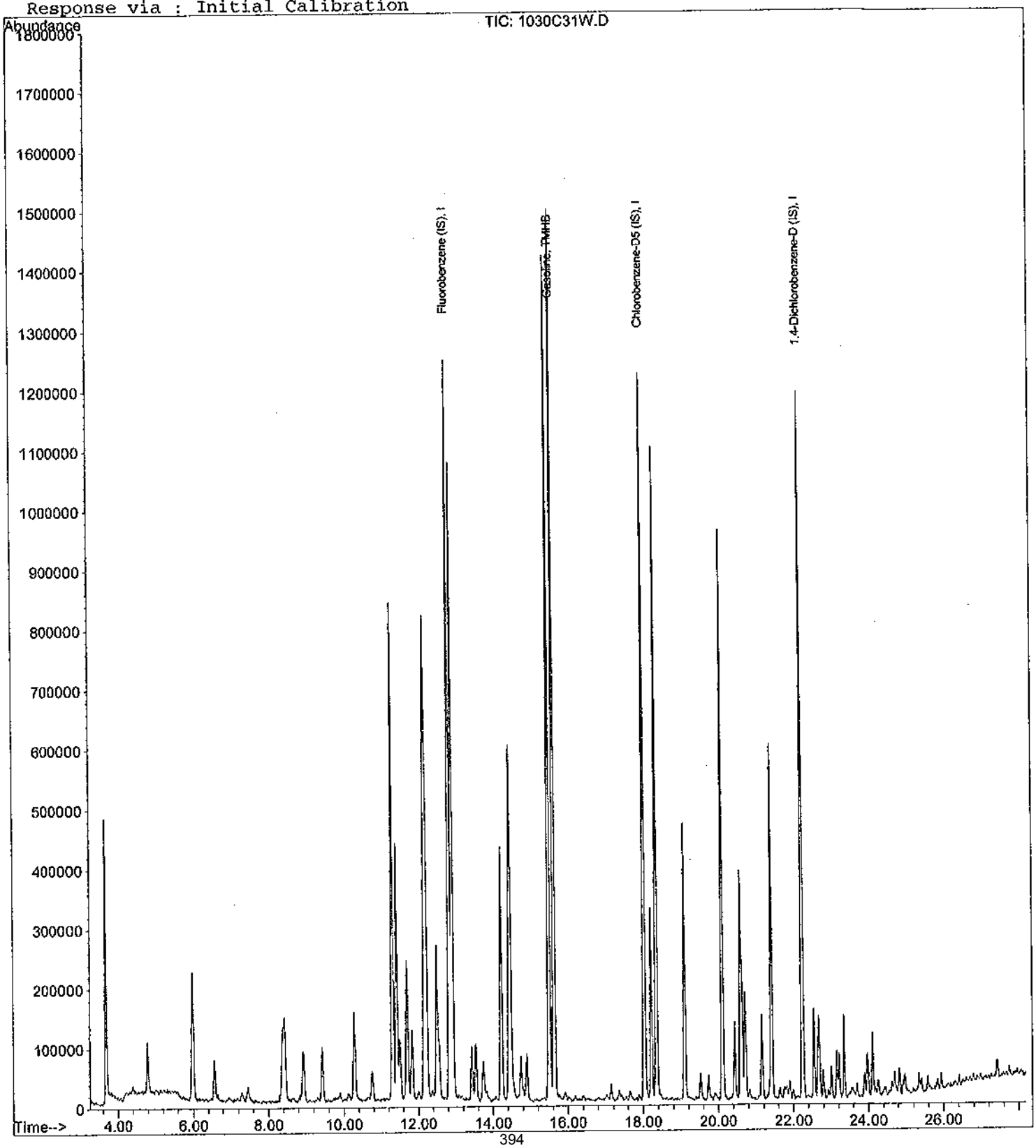
Quantitation Report

Data File : M:\CHICO\DATA\C111030\1030C31W.D
Acq On : 31 Oct 11 10:57
Sample : GAS 300ug/L LCS-1WC
Misc : Water 10mLw/ IS&S:10-30/10-26-11

Vial: 1
Operator: STC
Inst : Chico
Multiplr: 1.00

Quant Time: Nov 3 10:53 2011 Quant Results File: CGAS.RES

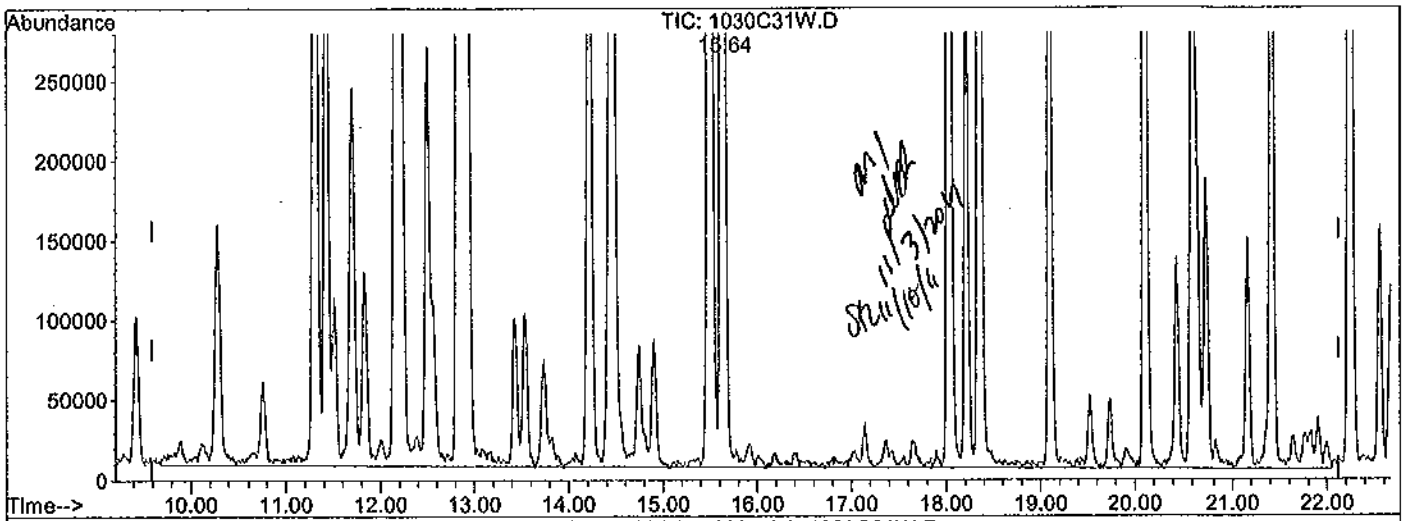
Method : M:\CHICO\DATA\C111030\CGAS.M (RTE Integrator)
Title : METHOD 8260B
Last Update : Thu Nov 03 10:47:02 2011
Response via : Initial Calibration

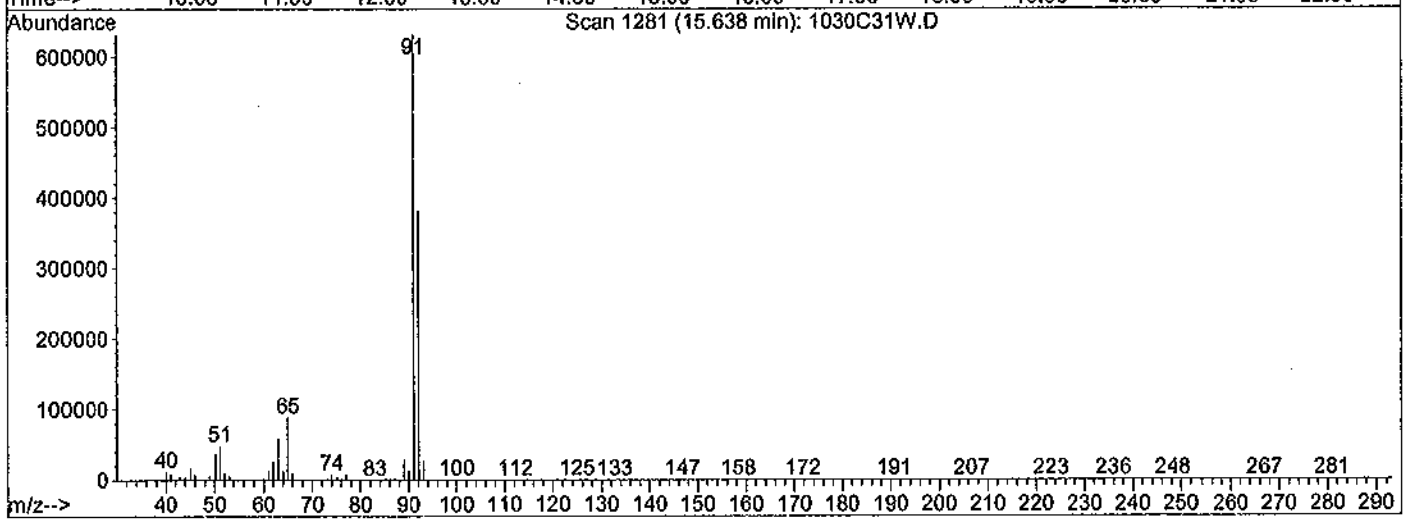


Quantitation Report

Data File : M:\CHICO\DATA\C111030\1030C31W.D Vial: 1
Acq On : 31 Oct 11 10:57 Operator: STC
Sample : GAA 300ug/L LCS-1WC Inst : Chico
Misc : Water 10mLw/ IS&S:10-30/10-26-11 Multiplr: 1.00
Quant Time: Nov 3 10:52 2011 Quant Results File: temp.res

Method : M:\CHICO\DATA\C111030\CGAS.M (RTE Integrator)
Title : METHOD 8260B
Last Update : Thu Nov 03 10:47:02 2011
Response via : Multiple Level Calibration





TIC: 1030C31W.D

(2) Gasoline (TMHB)

15.64min 380.8950ppb m

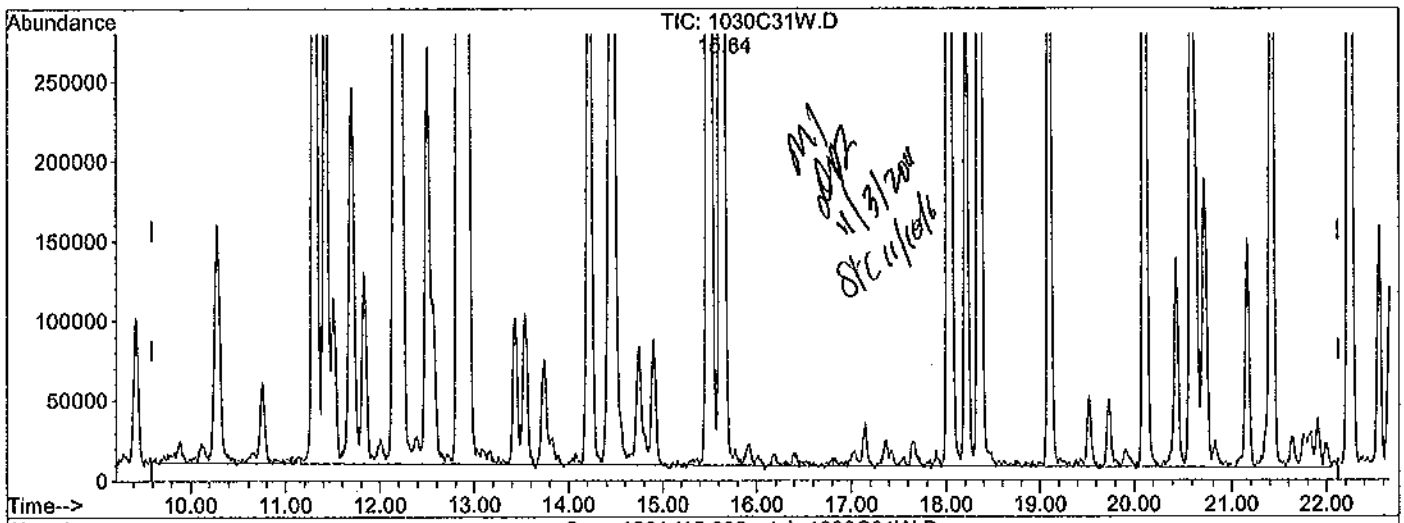
response 52602858

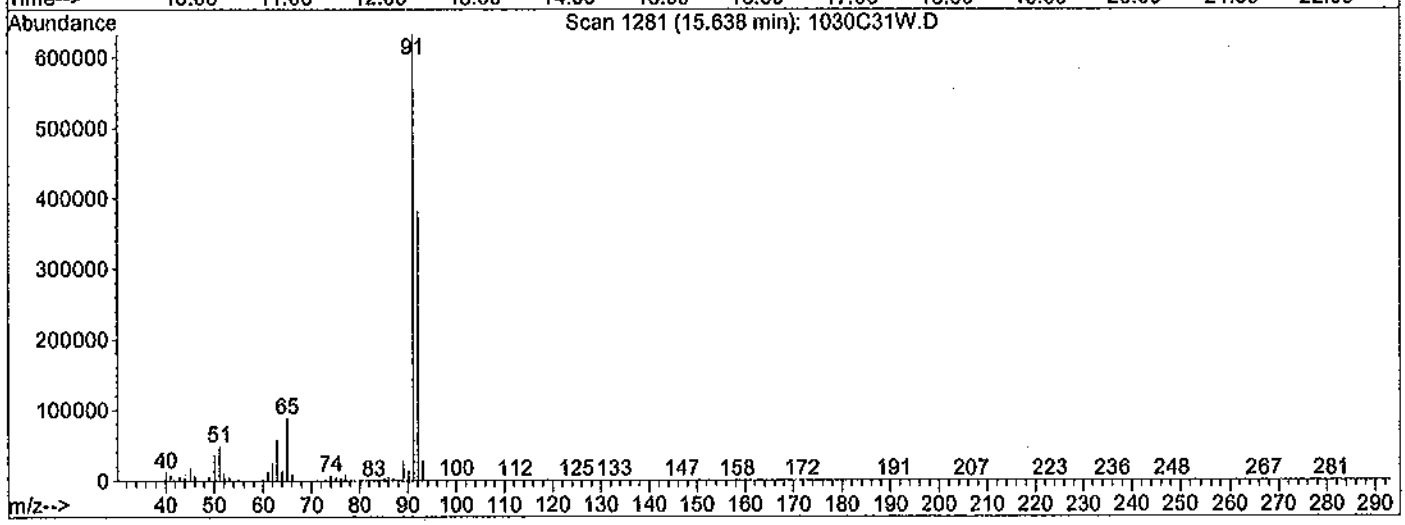
Ion	Exp%	Act%
TIC	100	100
0.00	0.00	0.21#
0.00	0.00	0.59#
0.00	0.00	0.00

Quantitation Report

Data File : M:\CHICO\DATA\C111030\1030C31W.D Vial: 1
Acq On : 31 Oct 11 10:57 Operator: STC
Sample : GAA 300ug/L LCS-1WC Inst : Chico
Misc : Water 10mLw/ IS&S:10-30/10-26-11 Multiplr: 1.00
Quant Time: Nov 3 10:53 2011 Quant Results File: temp.res

Method : M:\CHICO\DATA\C111030\CGAS.M (RTE Integrator)
Title : METHOD 8260B
Last Update : Thu Nov 03 10:47:02 2011
Response via : Multiple Level Calibration





TIC: 1030C31W.D

(2) Gasoline (TMHB)

15.64min 369.8272ppb m

response 51531735

Ion	Exp%	Act%
TIC	100	100
0.00	0.00	0.21#
0.00	0.00	0.60#
0.00	0.00	0.00

Matrix Spike Recoveries

EPA 8260B VOCs + Gas Water

APPL ID: **111031W-49334 MS - 161029**
Batch ID: #86RHB-111030AC
Sample ID: AY49334
Client ID: ES047

APPL Inc.
908 North Temperance Avenue
Clovis, CA 93611

Compound Name	Spike Lvl ug/L	Matrix Result ug/L	SPK Result ug/L	DUP Result ug/L	SPK % Recovery	DUP % Recovery	Recovery Limits	RPD %	RPD Limits
1,1,1,2-TETRACHLOROETHANE	10.00	ND	10.9	11.2	109	112	80-130	2.7	30
1,1,1-TRICHLOROETHANE	10.00	ND	8.95	9.51	89.5	95.1	65-130	6.1	30
1,1,2,2-TETRACHLOROETHANE	10.00	ND	12.7	12.6	127	126	65-130	0.79	30
1,1,2-TRICHLOROETHANE	10.00	ND	10.6	10.6	106	106	75-125	0.0	30
1,1-DICHLOROETHANE	10.00	ND	9.46	9.94	94.6	99.4	70-135	4.9	30
1,1-DICHLOROETHENE	10.00	ND	8.96	9.50	89.6	95.0	70-130	5.9	30
1,2,3-TRICHLOROPROPANE	10.00	ND	11.6	11.4	116	114	75-125	1.7	30
1,2,4-TRICHLOROBENZENE	10.00	ND	9.55	10.1	95.5	101	65-135	5.6	30
1,2-DIBROMO-3-CHLOROPROPANE	10.00	ND	11.2	12.2	112	122	50-130	8.5	30
1,2-DIBROMOETHANE	10.00	ND	10.9	10.8	109	108	70-130	0.92	30
1,2-DICHLOROBENZENE	10.00	ND	9.85	10.2	98.5	102	70-120	3.5	30
1,2-DICHLOROETHANE	10.00	ND	8.97	9.48	89.7	94.8	70-130	5.5	30
1,2-DICHLOROPROPANE	10.00	ND	9.62	10.5	96.2	105	75-125	8.7	30
1,3-DICHLOROBENZENE	10.00	ND	9.68	10.1	96.8	101	75-125	4.2	30
1,3-DICHLOROPROPENE, TOTAL	20.0	ND	20.4	20.9	102	105	70-130	2.4	30
1,4-DICHLOROBENZENE	10.00	ND	9.80	10.2	98.0	102	75-125	4.0	30
2-BUTANONE	10.00	ND	10.5	10.2	105	102	30-150	2.9	30
4-METHYL-2-PENTANONE	10.00	ND	11.0	11.0	110	110	60-135	0.0	30
ACETONE	10.00	ND	11.9	13.6	119	136	40-140	13.3	30
BENZENE	10.00	ND	9.56	10.2	95.6	102	80-120	6.5	30
BROMODICHLOROMETHANE	10.00	ND	9.85	10.3	98.5	103	75-120	4.5	30
BROMOFORM	10.00	ND	9.59	10.1	95.9	101	70-130	5.2	30
BROMOMETHANE	10.00	ND	9.35	9.85	93.5	98.5	30-145	5.2	30
CARBON TETRACHLORIDE	10.00	ND	9.77	9.93	97.7	99.3	65-140	1.6	30
CHLOROBENZENE	10.00	ND	9.68	10.1	96.8	101	80-120	4.2	30

= Recovery is outside QC limits.

Comments:

Primary	SPK	DUP
Quant Method :	CALLW.M	CALLW.M
Extraction Date :	10/31/11	10/31/11
Analysis Date :	10/31/11	10/31/11
Instrument :	Chico	Chico
Run :	1030C39	1030C40
Initials :	ARS	

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APPL MSD SCII

Matrix Spike Recoveries

EPA 8260B VOCs + Gas Water

APPL ID: **111031W-49334 MS - 161029**
Batch ID: #86RHB-111030AC
Sample ID: AY49334
Client ID: ES047

APPL Inc.
908 North Temperance Avenue
Clovis, CA 93611

Compound Name	Spike Lvl ug/L	Matrix Result ug/L	SPK Result ug/L	DUP Result ug/L	SPK % Recovery	DUP % Recovery	Recovery Limits	RPD %	RPD Limits
CHLORODIBROMOMETHANE	10.00	ND	10.8	11.4	108	114	60-135	5.4	30
CHLOROETHANE	10.00	ND	8.88	9.26	88.8	92.6	60-135	4.2	30
CHLOROFORM	10.00	ND	9.21	9.65	92.1	96.5	65-135	4.7	30
CHLOROMETHANE	10.00	ND	8.69	9.03	86.9	90.3	40-125	3.8	30
CIS-1,2-DICHLOROETHENE	10.00	ND	9.19	9.57	91.9	95.7	70-125	4.1	30
ETHYLBENZENE	10.00	ND	9.73	10.3	97.3	103	75-125	5.7	30
GASOLINE	300	ND	374	377	125	126 #	75-125	0.80	30
HEXACHLOROBUTADIENE	10.00	ND	9.04	9.78	90.4	97.8	50-140	7.9	30
METHYL TERT-BUTYL ETHER	10.00	ND	10.1	10.2	101	102	65-125	0.99	30
METHYLENE CHLORIDE	10.00	ND	9.83	9.84	98.3	98.4	55-140	0.10	30
STYRENE	10.00	ND	10.2	10.6	102	106	65-135	3.8	30
TETRACHLOROETHENE	10.00	ND	9.48	10.3	94.8	103	45-150	8.3	30
TOLUENE	10.00	ND	9.46	10.4	94.6	104	75-120	9.5	30
TRANS-1,2-DICHLOROETHENE	10.00	ND	9.17	9.47	91.7	94.7	60-140	3.2	30
TRICHLOROETHENE	10.00	ND	9.00	9.38	90.0	93.8	70-125	4.1	30
VINYL CHLORIDE	10.00	ND	9.90	9.78	99.0	97.8	50-145	1.2	30
XYLENES (TOTAL)	30.0	ND	29.1	30.7	97.0	102	80-120	5.4	30
SURROGATE: 1,2-DICHLOROETHANE-D	24.2	NA	22.7	22.3	93.7	92.1	70-120		
SURROGATE: 4-BROMOFLUOROBENZE	25.5	NA	25.8	26.7	101	105	75-120		
SURROGATE: DIBROMOFLUOROMETH	25.1	NA	24.7	25.0	98.4	99.6	85-115		
SURROGATE: TOLUENE-D8 (S)	25.8	NA	26.3	26.9	102	104	85-120		

= Recovery is outside QC limits.

Comments:

Primary	SPK	DUP
Quant Method :	CALLW.M	CALLW.M
Extraction Date :	10/31/11	10/31/11
Analysis Date :	10/31/11	10/31/11
Instrument :	Chico	Chico
Run :	1030C39	1030C40
Initials :	ARS	

Printed: 12/06/11 3:06:24 PM
APPL MSD SCII

Quantitation Report (Not Reviewed)

Data File : M:\CHICO\DATA\C111030\1030C39W.D Vial: 1
Acq On : 31 Oct 11 16:08 Operator: STC
Sample : AY49334W141516 MS-1WC (VOC) Inst : Chico
Misc : Water 10mLw/ IS&S:10-30/10-26-11 Multiplr: 1.00

Quant Time: Dec 2 11:35 2011 Quant Results File: CALLW.RES

Quant Method : M:\CHICO\DATA\C111030\CALLW.M (RTE Integrator)
Title : METHOD 8260
Last Update : Fri Dec 02 11:32:50 2011
Response via : Initial Calibration
DataAcq Meth : V8260

Internal Standards	R.T.	QIon	Response	Conc Units		Dev(Min)
1) Fluorobenzene (IS)	12.85	96	673315	25.00000 ppb		0.02
55) Chlorobenzene-D5 (IS)	18.05	117	444608	25.00000 ppb		0.00
71) 1,4-Dichlorobenzene-D (IS)	22.26	152	234048	25.00000 ppb		0.02
System Monitoring Compounds						
33) Dibromofluoromethane(S)	11.43	111	443656	24.73529 ppb		0.00
Spiked Amount 25.097			Recovery = 98.557%			
38) 1,2-DCA-D4(S)	12.23	65	361670	22.65208 ppb		0.00
Spiked Amount 24.225			Recovery = 93.506%			
56) Toluene-D8(S)	15.51	98	1644192	26.28251 ppb		0.02
Spiked Amount 25.808			Recovery = 101.839%			
64) 4-Bromofluorobenzene(S)	20.12	95	578226	25.79618 ppb		0.00
Spiked Amount 25.459			Recovery = 101.322%			
Target Compounds						Qvalue
2) Dichlorodifluoromethane	4.08	85	233739	9.42228 ppb		96
3) Freon 114	4.34	85	156235	10.05590 ppb		100
4) Chloromethane	4.56	50	267000	8.68546 ppb		98
5) Vinyl chloride	4.82	62	203545	9.89891 ppb		95
7) Bromomethane	5.73	94	139543	9.34878 ppb		90
8) Chloroethane	5.92	64	150908	8.88492 ppb		95
9) Dichlorofluoromethane	6.01	67	417786	8.89643 ppb		98
10) Trichlorofluoromethane	6.55	101	259316	9.30510 ppb		97
11) Acetonitrile	7.67	41	102299	138.69122 ug/l		100
12) Acrolein	7.18	56	41239	122.17811 ppb		98
13) Acetone	7.30	43	23028	11.93711 ppb	#	67
14) Freon-113	7.48	101	166680	10.24118 ppb		91
15) 1,1-DCE	7.69	96	172282	8.96239 ppb		93
16) t-Butanol	7.78	59	13291	145.60910 ppb		93
17) Methyl Acetate	8.21	43	57808	9.94422 ppb		98
18) Iodomethane	8.17	142	85074	8.93041 ppb		91
19) Acrylonitrile	8.59	53	23274	11.05716 ppb		92
20) Methylene chloride	8.48	84	180285	9.83293 ppb		97
21) Carbon disulfide	8.57	76	184128	9.85829 ppb		99
22) Methyl t-butyl ether (MtBE	8.92	73	294046	10.12039 ppb		91
23) Trans-1,2-DCE	9.11	96	204403	9.16612 ppb		91
24) Diisopropyl Ether	9.77	45	657158	10.23091 ppb		91
25) 1,1-DCA	9.80	63	360267	9.45758 ppb		96
26) Vinyl Acetate	9.43	43	123424	10.42529 ppb		85
27) Ethyl tert Butyl Ether	10.46	59	458722	10.46060 ppb		99
28) MEK (2-Butanone)	10.44	43	79943	10.46895 ppb		94
29) Cis-1,2-DCE	10.83	96	210715	9.19494 ppb		93
30) 2,2-Dichloropropane	10.82	77	272392	9.98243 ppb		100
31) Chloroform	11.10	83	337366	9.20607 ppb		100
32) Bromochloromethane	11.33	128	66189	10.37457 ppb	#	71
34) 1,1,1-TCA	11.85	97	298219	8.94840 ppb		96
35) Cyclohexane	12.01	56	300163	9.67354 ppb		99
36) 1,1-Dichloropropene	12.12	75	263597	9.23085 ppb		96
37) 2,2,4-Trimethylpentane	12.19	57	501585	10.31895 ppb		95
39) Carbon Tetrachloride	12.31	117	224210	9.77018 ppb		100
40) Tert Amyl Methyl Ether	12.36	73	337649	10.29749 ppb	#	95

(#) = qualifier out of range (m) = manual integration

Quantitation Report (Not Reviewed)

Data File : M:\CHICO\DATA\C111030\1030C39W.D Vial: 1
Acq On : 31 Oct 11 16:08 Operator: STC
Sample : AY49334W141516 MS-1WC (VOC) Inst : Chico
Misc : Water 10mLw/ IS&S:10-30/10-26-11 Multiplr: 1.00

Quant Time: Dec 2 11:35 2011 Quant Results File: CALLW.RES

Quant Method : M:\CHICO\DATA\C111030\CALLW.M (RTE Integrator)
Title : METHOD 8260
Last Update : Fri Dec 02 11:32:50 2011
Response via : Initial Calibration
DataAcq Meth : V8260

	Compound	R.T.	QIon	Response	Conc	Unit		Qvalue
41)	1,2-DCA	12.39	62	168650	8.96568	ppb		99
42)	Benzene	12.52	78	784311	9.55972	ppb		98
43)	TCE	13.55	95	204442	8.99750	ppb		90
44)	2-Pentanone	13.21	43	641372	134.94011	ppb		98
45)	1,2-Dichloropropane	13.77	63	179306	9.61525	ppb	#	92
46)	Bromodichloromethane	14.13	83	209814	9.84828	ppb		85
47)	Methyl Cyclohexane	13.83	83	264020	9.94314	ppb		99
48)	Dibromomethane	14.18	93	75188	10.08162	ppb		96
49)	2-Chloroethyl vinyl ether	14.59	63	44799	9.44967	ppb		97
50)	1-Bromo-2-chloroethane	14.89	63	159282	10.01037	ppb		91
51)	Cis-1,3-Dichloropropene	15.02	75	212722	10.47106	ppb		97
52)	Toluene	15.65	91	765314	9.45759	ppb		97
53)	Trans-1,3-Dichloropropene	15.81	75	145848	9.97216	ppb		99
54)	1,1,2-TCA	16.09	83	83942	10.64886	ppb		94
57)	1,2-EDB	17.34	107	92334	10.91173	ppb		92
58)	Tetrachloroethene	16.80	164	216714	9.48077	ppb		97
59)	1-Chlorohexane	17.72	91	263885	10.02502	ppb		88
60)	1,1,1,2-Tetrachloroethane	18.17	131	156166	10.91277	ppb		96
61)	m&p-Xylene	18.37	106	646753	19.15250	ppb		99
62)	o-Xylene	19.11	106	322766	9.93837	ppb		99
63)	Styrene	19.13	104	500812	10.21655	ppb		96
65)	2-Hexanone	16.11	43	49009	12.04661	ppb		98
66)	1,3-Dichloropropane	16.51	76	172305	10.32618	ppb		99
67)	Dibromochloromethane	16.98	129	117679	10.80307	ppb		87
68)	Chlorobenzene	18.12	112	467466	9.67690	ppb		98
69)	Ethylbenzene	18.23	91	875108	9.72830	ppb		98
70)	Bromoform	19.65	173	53217	9.59314	ppb		95
72)	MIBK (methyl isobutyl keto	14.68	43	73055	10.97983	ppb		80
73)	Isopropylbenzene	19.75	105	1154759	13.61308	ppb		99
74)	1,1,2,2-Tetrachloroethane	19.92	83	90087	12.68678	ppb	#	76
75)	1,2,3-Trichloropropane	20.17	110	9227	11.57218	ppb		78
76)	t-1,4-Dichloro-2-Butene	20.23	53	20628	12.81166	ppb	#	78
77)	Bromobenzene	20.49	156	193805	9.90473	ppb		92
78)	n-Propylbenzene	20.46	91	1397086	13.80582	ppb		98
79)	4-Ethyltoluene	20.65	105	653523	9.33219	ppb		98
80)	2-Chlorotoluene	20.75	91	606805	9.05395	ppb		100
81)	1,3,5-Trimethylbenzene	20.73	105	642904	9.33159	ppb		98
82)	4-Chlorotoluene	20.82	91	563219	9.75925	ppb		98
83)	Tert-Butylbenzene	21.37	119	771335	10.34106	ppb		98
84)	1,2,4-Trimethylbenzene	21.43	105	651620	9.05592	ppb		97
85)	Sec-Butylbenzene	21.77	105	1244018	13.90727	ppb		100
86)	p-Isopropyltoluene	22.00	119	725182	9.46493	ppb		99
87)	Benzyl Chloride	22.44	91	197410	19.40978	ppb		97
88)	1,3-DCB	22.14	146	387323	9.67943	ppb		96
89)	1,4-DCB	22.31	146	363787	9.79512	ppb		92
90)	Hexachloroethane	23.56	117	533859	38.29226	ppb	#	12
91)	n-Butylbenzene	22.71	91	769919	11.52107	ppb		98
92)	1,2-DCB	22.93	146	313462	9.84841	ppb		98
93)	1,2-Dibromo-3-chloropropan	24.14	155	12899	11.18801	ppb		82
94)	1,2,4-Trichlorobenzene	25.59	180	220157	9.54544	ppb		98
95)	Hexachlorobutadiene	25.86	223	37896	9.04368	ppb		97

(#) = qualifier out of range (m) = manual integration

Quantitation Report (Not Reviewed)

Data File : M:\CHICO\DATA\C111030\1030C39W.D Vial: 1
Acq On : 31 Oct 11 16:08 Operator: STC
Sample : AY49334W141516 MS-1WC (VOC) Inst : Chico
Misc : Water 10mLw/ IS&S:10-30/10-26-11 Multiplr: 1.00

Quant Time: Dec 2 11:35 2011 Quant Results File: CALLW.RES

Quant Method : M:\CHICO\DATA\C111030\CALLW.M (RTE Integrator)
Title : METHOD 8260
Last Update : Fri Dec 02 11:32:50 2011
Response via : Initial Calibration
DataAcq Meth : V8260

Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
96) Naphthalene	25.94	128	595126	20.91304	ppb	100
97) 1,2,3-Trichlorobenzene	26.31	180	169004	9.68724	ppb	98

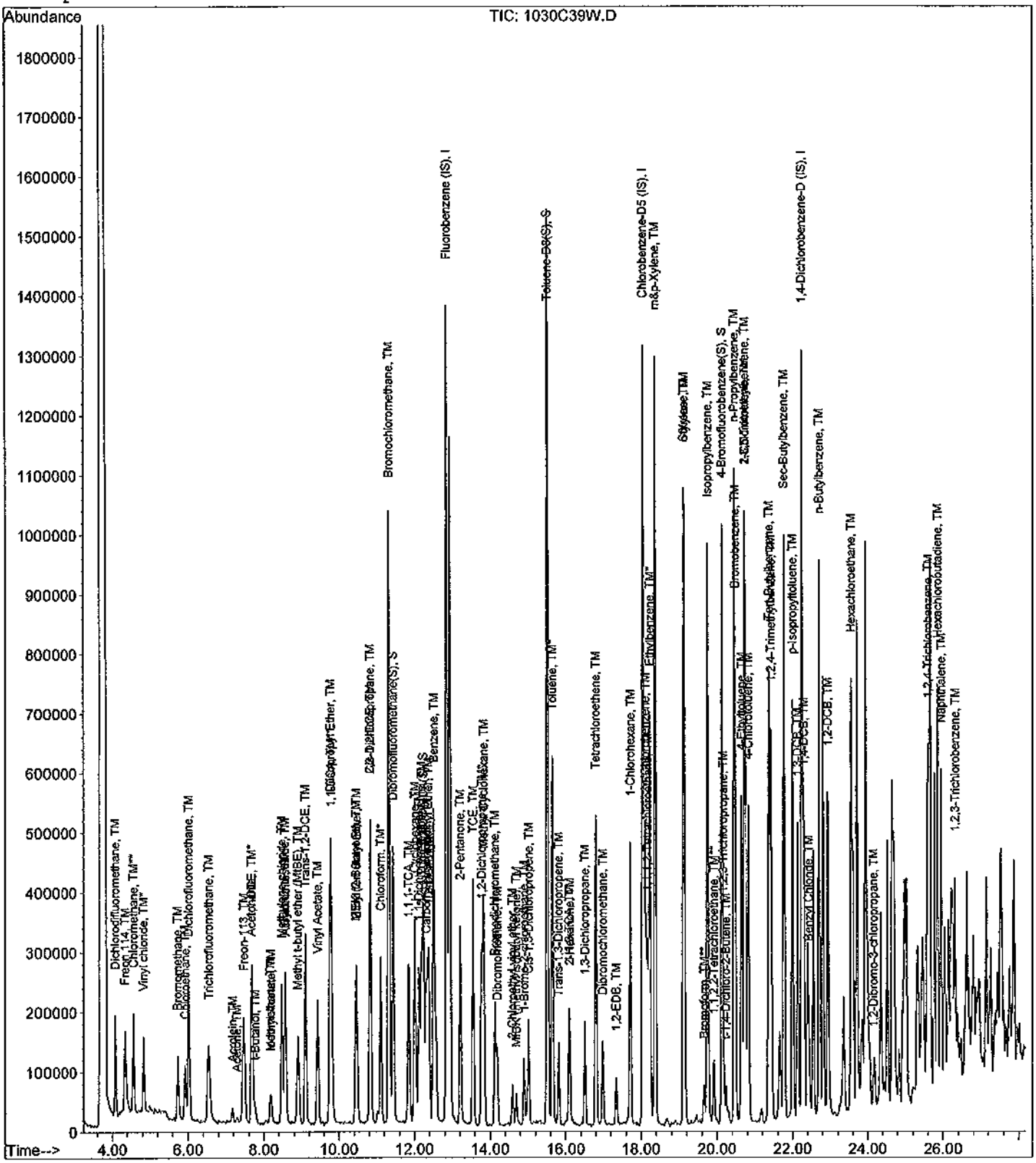
(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : M:\CHICO\DATA\C111030\1030C39W.D Vial: 1
Acq On : 31 Oct 11 16:08 Operator: STC
Sample : AY49334W141516 MS-1WC (VOC) Inst : Chico
Misc : Water 10mLw/ IS&S:10-30/10-26-11 Multiplr: 1.00

Quant Time: Dec 2 11:35 2011 Quant Results File: CALLW.RES

Method : M:\CHICO\DATA\C111030\CALLW.M (RTE Integrator)
Title : METHOD 8260
Last Update : Fri Dec 02 11:32:50 2011
Response via : Initial Calibration



Quantitation Report (Not Reviewed)

Data File : M:\CHICO\DATA\C111030\1030C40W.D Vial: 1
Acq On : 31 Oct 11 16:45 Operator: STC
Sample : AY49334W141516 MSD-1WC (VOC) Inst : Chico
Misc : Water 10mLw/ IS&S:10-30/10-26-11 Multiplr: 1.00

Quant Time: Dec 2 11:35 2011 Quant Results File: CALLW.RES

Quant Method : M:\CHICO\DATA\C111030\CALLW.M (RTE Integrator)
Title : METHOD 8260
Last Update : Fri Dec 02 11:32:50 2011
Response via : Initial Calibration
DataAcq Meth : V8260

Internal Standards	R.T.	QIon	Response	Conc	Units		Dev(Min)
1) Fluorobenzene (IS)	12.85	96	673885	25.00000	ppb		0.02
55) Chlorobenzene-D5 (IS)	18.05	117	449984	25.00000	ppb		0.00
71) 1,4-Dichlorobenzene-D (IS)	22.26	152	240256	25.00000	ppb		0.02
System Monitoring Compounds							
33) Dibromofluoromethane(S)	11.43	111	449233	25.02504	ppb		0.00
Spiked Amount 25.097			Recovery = 99.712%				
38) 1,2-DCA-D4(S)	12.23	65	356030	22.27998	ppb		0.00
Spiked Amount 24.225			Recovery = 91.970%				
56) Toluene-D8(S)	15.51	98	1704225	26.91667	ppb		0.02
Spiked Amount 25.808			Recovery = 104.296%				
64) 4-Bromofluorobenzene(S)	20.13	95	605686	26.69841	ppb		0.02
Spiked Amount 25.459			Recovery = 104.865%				
Target Compounds							Qvalue
2) Dichlorodifluoromethane	4.08	85	220598	8.88503	ppb		94
3) Freon 114	4.34	85	150416	9.67318	ppb		95
4) Chloromethane	4.56	50	277770	9.02816	ppb		93
5) Vinyl chloride	4.82	62	201338	9.78330	ppb		96
7) Bromomethane	5.74	94	147121	9.84814	ppb		97
8) Chloroethane	5.93	64	157372	9.25766	ppb		99
9) Dichlorofluoromethane	6.01	67	450238	9.57936	ppb		98
10) Trichlorofluoromethane	6.54	101	257325	9.22585	ppb		95
11) Acetonitrile	7.66	41	109689	148.58438	ug/l		100
12) Acrolein	7.18	56	40703	120.48812	ppb		94
13) Acetone	7.29	43	26277	13.60978	ppb	#	76
14) Freon-113	7.48	101	164332	10.07108	ppb		98
15) 1,1-DCE	7.70	96	182764	9.49964	ppb		92
16) t-Butanol	7.78	59	14747	161.42360	ppb		98
17) Methyl Acetate	8.20	43	62200	10.74239	ppb		99
18) Iodomethane	8.18	142	105131	10.12939	ppb		89
19) Acrylonitrile	8.58	53	23406	11.11230	ppb		88
20) Methylene chloride	8.48	84	180541	9.83857	ppb		95
21) Carbon disulfide	8.58	76	183104	9.79517	ppb		100
22) Methyl t-butyl ether (MtBE	8.92	73	297031	10.21448	ppb		94
23) Trans-1,2-DCE	9.11	96	211447	9.47397	ppb		98
24) Diisopropyl Ether	9.77	45	672476	10.46053	ppb		91
25) 1,1-DCA	9.80	63	378977	9.94033	ppb		97
26) Vinyl Acetate	9.44	43	132519	11.29670	ppb		98
27) Ethyl tert Butyl Ether	10.45	59	467350	10.64834	ppb		99
28) MEK (2-Butanone)	10.44	43	78228	10.22050	ppb		95
29) Cis-1,2-DCE	10.82	96	219402	9.56592	ppb		95
30) 2,2-Dichloropropane	10.82	77	284485	10.41678	ppb		99
31) Chloroform	11.10	83	353899	9.64906	ppb		100
32) Bromochloromethane	11.33	128	69588	10.89810	ppb		79
34) 1,1,1-TCA	11.85	97	317360	9.51469	ppb		95
35) Cyclohexane	12.01	56	307089	9.88838	ppb		94
36) 1,1-Dichloropropene	12.12	75	277564	9.71173	ppb		94
37) 2,2,4-Trimethylpentane	12.20	57	511508	10.53601	ppb		91
39) Carbon Tetrachloride	12.31	117	228003	9.92706	ppb		94
40) Tert Amyl Methyl Ether	12.36	73	345658	10.53283	ppb		97

(#) = qualifier out of range (m) = manual integration

Quantitation Report (Not Reviewed)

Data File : M:\CHICO\DATA\C111030\1030C40W.D Vial: 1
Acq On : 31 Oct 11 16:45 Operator: STC
Sample : AY49334W141516 MSD-1WC (VOC) Inst : Chico
Misc : Water 10mLw/ IS&S:10-30/10-26-11 Multiplr: 1.00

Quant Time: Dec 2 11:35 2011 Quant Results File: CALLW.RES

Quant Method : M:\CHICO\DATA\C111030\CALLW.M (RTE Integrator)
Title : METHOD 8260
Last Update : Fri Dec 02 11:32:50 2011
Response via : Initial Calibration
DataAcq Meth : V8260

Compound	R.T.	QIon	Response	Conc	Unit		Qvalue
41) 1,2-DCA	12.39	62	178492	9.48087	ppb		94
42) Benzene	12.52	78	833883	10.15534	ppb		99
43) TCE	13.55	95	213329	9.38068	ppb		92
44) 2-Pentanone	13.21	43	670030	140.85032	ppb		94
45) 1,2-Dichloropropane	13.77	63	195266	10.46224	ppb		95
46) Bromodichloromethane	14.13	83	219583	10.29810	ppb		94
47) Methyl Cyclohexane	13.83	83	263113	9.90060	ppb		97
48) Dibromomethane	14.19	93	77194	10.34184	ppb		90
49) 2-Chloroethyl vinyl ether	14.59	63	43646	9.19868	ppb	#	93
50) 1-Bromo-2-chloroethane	14.89	63	166408	10.44937	ppb		90
51) Cis-1,3-Dichloropropene	15.02	75	211155	10.38513	ppb		96
52) Toluene	15.65	91	840102	10.37302	ppb		98
53) Trans-1,3-Dichloropropene	15.81	75	154386	10.54700	ppb		98
54) 1,1,2-TCA	16.09	83	83579	10.59385	ppb	#	88
57) 1,2-EDB	17.34	107	92512	10.80216	ppb		95
58) Tetrachloroethene	16.80	164	239347	10.34582	ppb		93
59) 1-Chlorohexane	17.72	91	280271	10.52032	ppb		94
60) 1,1,1,2-Tetrachloroethane	18.17	131	161501	11.15075	ppb		87
61) m&p-Xylene	18.37	106	683414	19.99636	ppb		98
62) o-Xylene	19.11	106	353111	10.74283	ppb		99
63) Styrene	19.13	104	527765	10.63776	ppb		99
65) 2-Hexanone	16.12	43	51343	12.46954	ppb		89
66) 1,3-Dichloropropane	16.51	76	177186	10.49183	ppb		100
67) Dibromochloromethane	16.98	129	125506	11.38395	ppb		93
68) Chlorobenzene	18.12	112	491969	10.06246	ppb		98
69) Ethylbenzene	18.23	91	936335	10.28458	ppb		100
70) Bromoform	19.64	173	56899	10.07037	ppb		90
72) MIBK (methyl isobutyl keto	14.69	43	75276	11.02130	ppb	#	77
73) Isopropylbenzene	19.75	105	1237307	14.20931	ppb		99
74) 1,1,2,2-Tetrachloroethane	19.91	83	91721	12.58313	ppb	#	89
75) 1,2,3-Trichloropropane	20.17	110	9365	11.44279	ppb	#	75
76) t-1,4-Dichloro-2-Butene	20.23	53	22200	13.43173	ppb		93
77) Bromobenzene	20.49	156	206516	10.28163	ppb		100
78) n-Propylbenzene	20.46	91	1530297	14.73145	ppb		98
79) 4-Ethyltoluene	20.65	105	706238	9.82436	ppb		95
80) 2-Chlorotoluene	20.75	91	653382	9.49701	ppb		97
81) 1,3,5-Trimethylbenzene	20.73	105	702310	9.93046	ppb		99
82) 4-Chlorotoluene	20.82	91	605738	10.22479	ppb		99
83) Tert-Butylbenzene	21.38	119	820988	10.72234	ppb		97
84) 1,2,4-Trimethylbenzene	21.43	105	691895	9.36719	ppb		99
85) Sec-Butylbenzene	21.77	105	1352492	14.72924	ppb		98
86) p-Isopropyltoluene	22.00	119	794796	10.10548	ppb		99
87) Benzyl Chloride	22.44	91	198774	19.03890	ppb		98
88) 1,3-DCB	22.14	146	415514	10.11563	ppb		95
89) 1,4-DCB	22.30	146	389740	10.22276	ppb		98
90) Hexachloroethane	23.56	117	551389	38.51564	ppb	#	12
91) n-Butylbenzene	22.71	91	850928	12.40428	ppb		97
92) 1,2-DCB	22.93	146	334523	10.23853	ppb		97
93) 1,2-Dibromo-3-chloropropan	24.14	155	14637	12.23153	ppb		89
94) 1,2,4-Trichlorobenzene	25.59	180	240070	10.13986	ppb		96
95) Hexachlorobutadiene	25.85	223	42056	9.77711	ppb		96

(#) = qualifier out of range (m) = manual integration

Quantitation Report (Not Reviewed)

Data File : M:\CHICO\DATA\C111030\1030C40W.D Vial: 1
Acq On : 31 Oct 11 16:45 Operator: STC
Sample : AY49334W141516 MSD-1WC (VOC) Inst : Chico
Misc : Water 10mLw/ IS&S:10-30/10-26-11 Multiplr: 1.00

Quant Time: Dec 2 11:35 2011 Quant Results File: CALLW.RES

Quant Method : M:\CHICO\DATA\C111030\CALLW.M (RTE Integrator)
Title : METHOD 8260
Last Update : Fri Dec 02 11:32:50 2011
Response via : Initial Calibration
DataAcq Meth : V8260

	Compound	R.T.	QIon	Response	Conc	Unit	Qvalue
96)	Naphthalene	25.94	128	619605	21.21064	ppb	99
97)	1,2,3-Trichlorobenzene	26.31	180	193601	10.81039	ppb	92

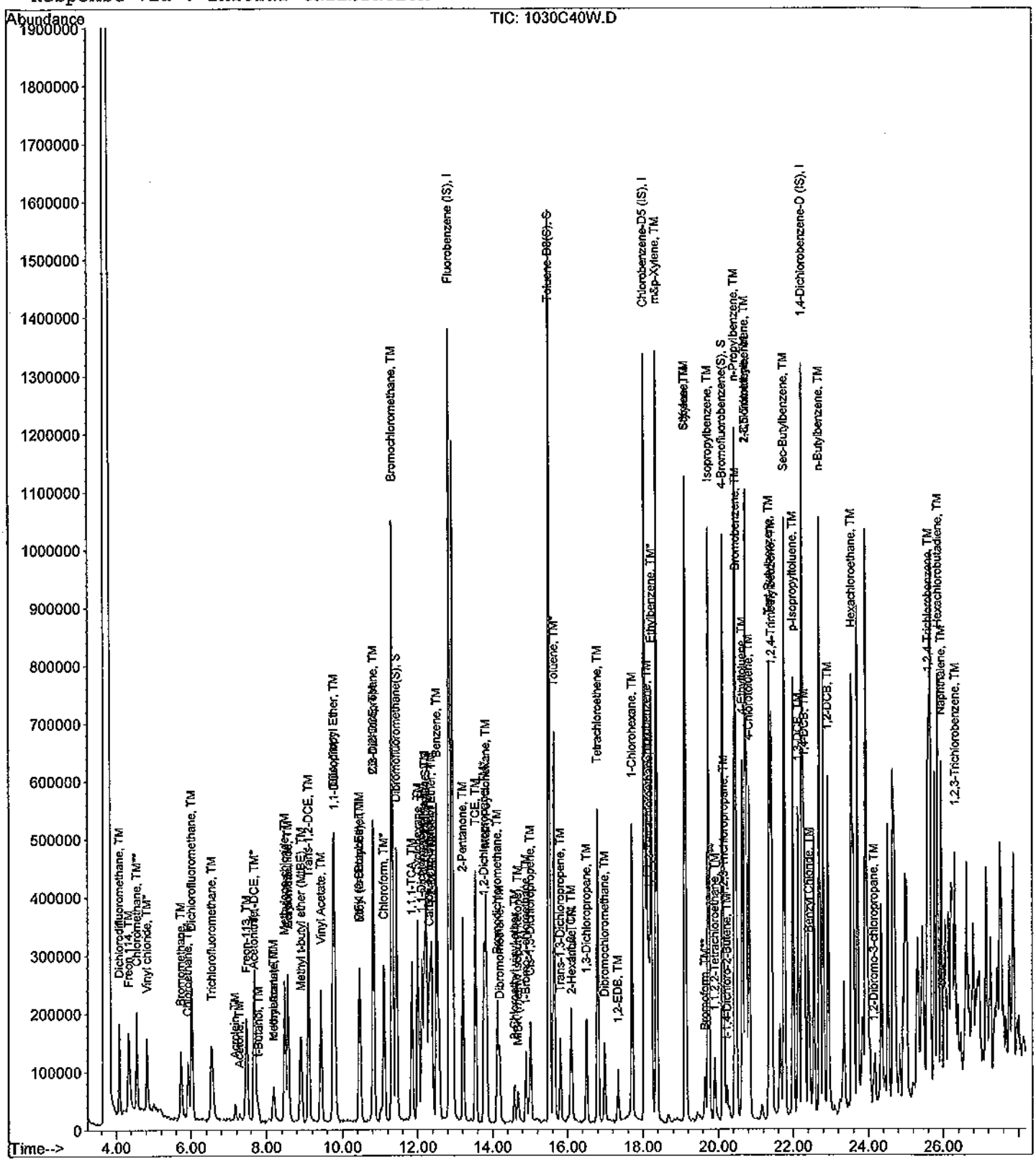
(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : M:\CHICO\DATA\C111030\1030C40W.D Vial: 1
Acq On : 31 Oct 11 16:45 Operator: STC
Sample : AY49334W141516 MSD-1WC (VOC) Inst : Chico
Misc : Water 10mLw/ IS&S:10-30/10-26-11 Multiplr: 1.00

Quant Time: Dec 2 11:35 2011 Quant Results File: CALLW.RES

Method : M:\CHICO\DATA\C111030\CALLW.M (RTE Integrator)
Title : METHOD 8260
Last Update : Fri Dec 02 11:32:50 2011
Response via : Initial Calibration



Quantitation Report (Not Reviewed)

Data File : M:\CHICO\DATA\C111030\1030C41W.D Vial: 1
Acq On : 31 Oct 11 17:22 Operator: STC
Sample : AY49334W161718 MS-1WC (GAS) Inst : Chico
Misc : Water 10mLw/ IS&S:10-30/10-26-11 Multiplr: 1.00

Quant Time: Nov 10 10:25 2011 Quant Results File: CGAS.RES

Quant Method : M:\CHICO\DATA\C111030\CGAS.M (RTE Integrator)
Title : METHOD 8260B
Last Update : Thu Nov 03 10:47:02 2011
Response via : Initial Calibration
DataAcq Meth : V8260

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Fluorobenzene (IS)	12.85	TIC	1408321	25.00000	ppb	0.01
3) Chlorobenzene-D5 (IS)	18.05	TIC	1375208	25.00000	ppb	0.01
4) 1,4-Dichlorobenzene-D (IS)	22.25	TIC	1432024	25.00000	ppb	0.01

System Monitoring Compounds

Target Compounds						Qvalue
2) Gasoline	15.64	TIC	59067164m	373.93954	ppb	100

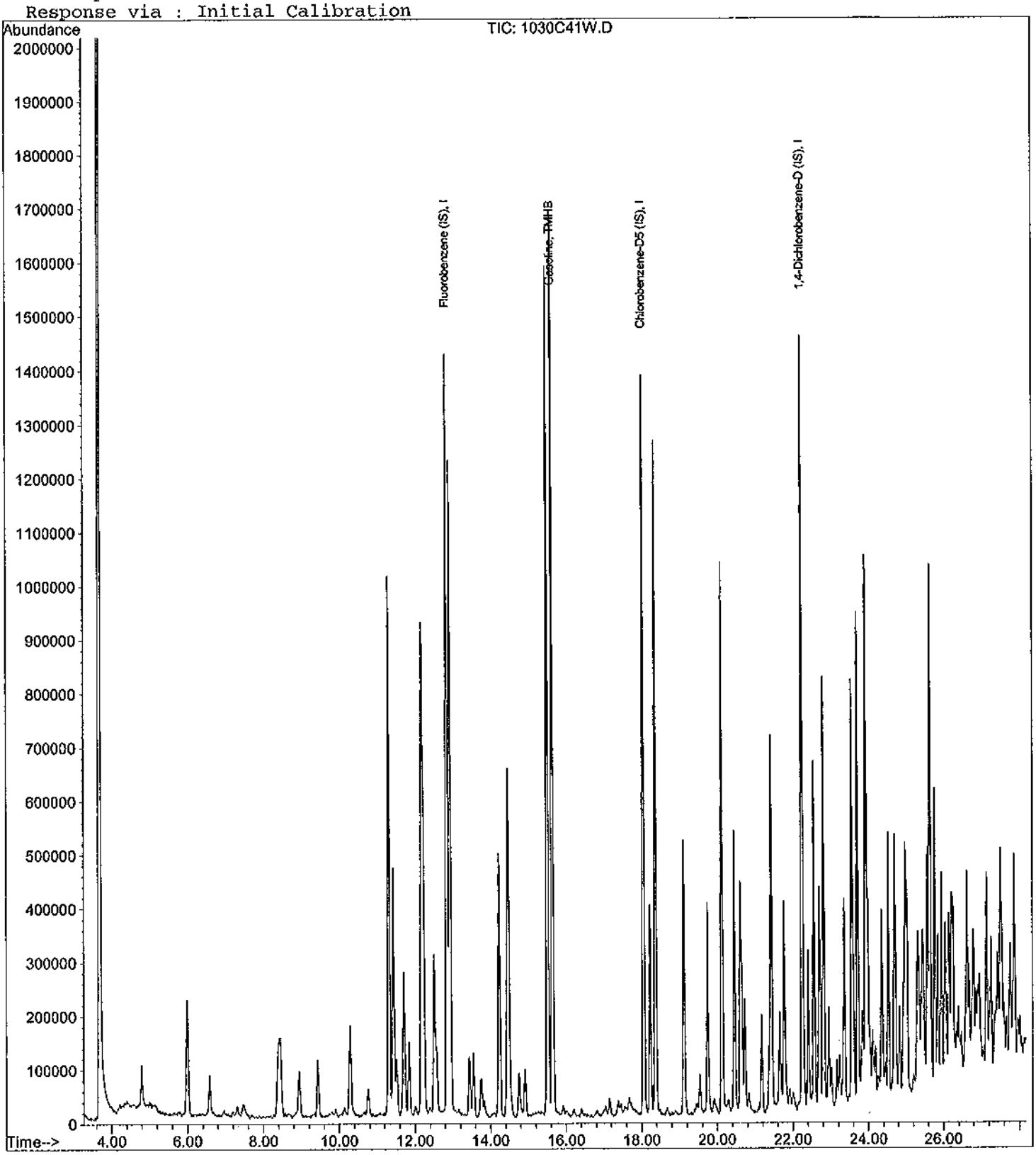
(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : M:\CHICO\DATA\C111030\1030C41W.D Vial: 1
Acq On : 31 Oct 11 17:22 Operator: STC
Sample : AY49334W161718 MS-1WC (GAS) Inst : Chico
Misc : Water 10mLw/ IS&S:10-30/10-26-11 Multiplr: 1.00

Quant Time: Nov 10 10:25 2011 Quant Results File: CGAS.RES

Method : M:\CHICO\DATA\C111030\CGAS.M (RTE Integrator)
Title : METHOD 8260B
Last Update : Thu Nov 03 10:47:02 2011
Response via : Initial Calibration



Quantitation Report (Not Reviewed)

Data File : M:\CHICO\DATA\C111030\1030C42W.D Vial: 1
Acq On : 31 Oct 11 17:59 Operator: STC
Sample : AY49334W161718 MSD-1WC (GAS) Inst : Chico
Misc : Water 10mLw/ IS&S:10-30/10-26-11 Multiplr: 1.00

Quant Time: Nov 3 10:23 2011 Quant Results File: CGAS.RES

Quant Method : M:\CHICO\DATA\C111030\CGAS.M (RTE Integrator)
Title : METHOD 8260B
Last Update : Thu Nov 03 10:47:02 2011
Response via : Initial Calibration
DataAcq Meth : V8260

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Fluorobenzene (IS)	12.85	TIC	1418194	25.00000	ppb	0.01
3) Chlorobenzene-D5 (IS)	18.05	TIC	1424242	25.00000	ppb	0.01
4) 1,4-Dichlorobenzene-D (IS)	22.26	TIC	1438638	25.00000	ppb	0.01

System Monitoring Compounds

Target Compounds						Qvalue
2) Gasoline	15.64	TIC	59788348m	376.70985	ppb	100

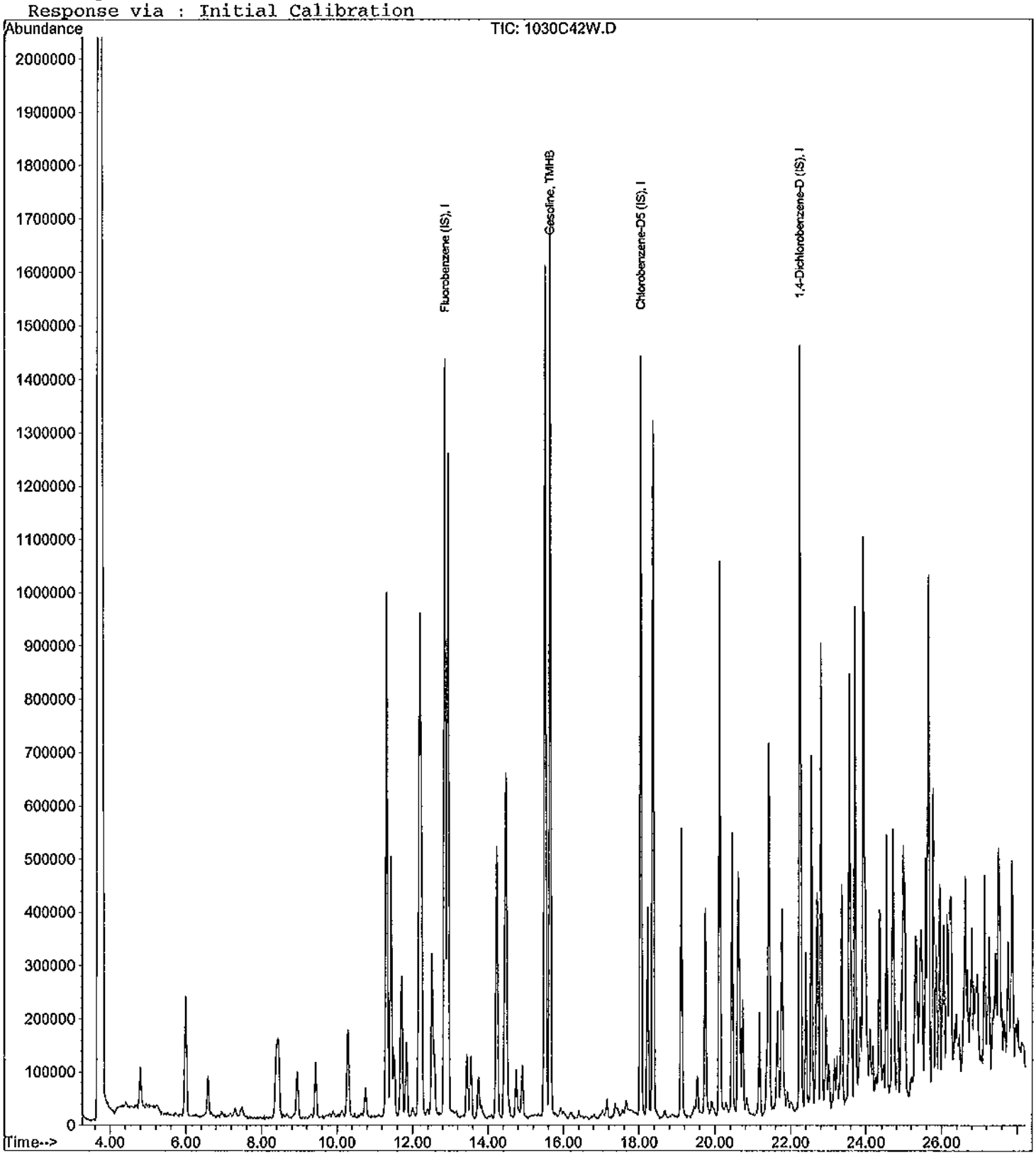
(#) = qualifier out of range (m) = manual integration

Quantitation Report

Data File : M:\CHICO\DATA\C111030\1030C42W.D Vial: 1
Acq On : 31 Oct 11 17:59 Operator: STC
Sample : AY49334W161718 MSD-1WC (GAS) Inst : Chico
Misc : Water 10mLw/ IS&S:10-30/10-26-11 Multiplr: 1.00

Quant Time: Nov 3 10:23 2011 Quant Results File: CGAS.RES

Method : M:\CHICO\DATA\C111030\CGAS.M (RTE Integrator)
Title : METHOD 8260B
Last Update : Thu Nov 03 10:47:02 2011
Response via : Initial Calibration

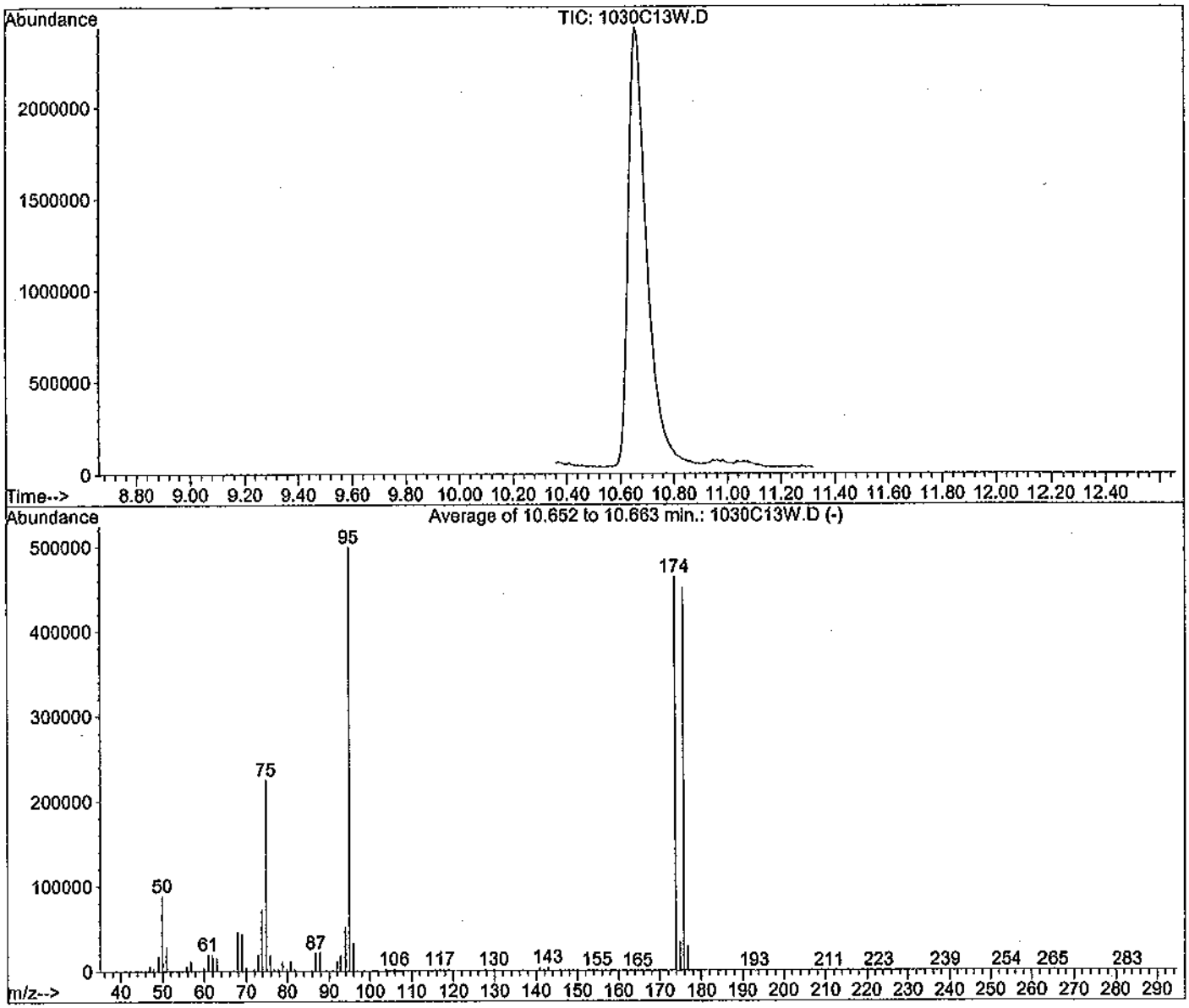


BFB

Data File : M:\CHICO\DATA\C111030\1030C13W.D
Acq On : 30 Oct 11 22:01
Sample : 20ug/ml BFB Std 10-19-11
Misc : Water 2uL

Vial: 1
Operator: RS
Inst : Chico
Multiplr: 1.00

Method : M:\CHICO\DATA\C111030\CALLW.M (RTE Integrator)
Title : METHOD 8260



AutoFind: Scans 52, 53, 54; Background Corrected with Scan 36

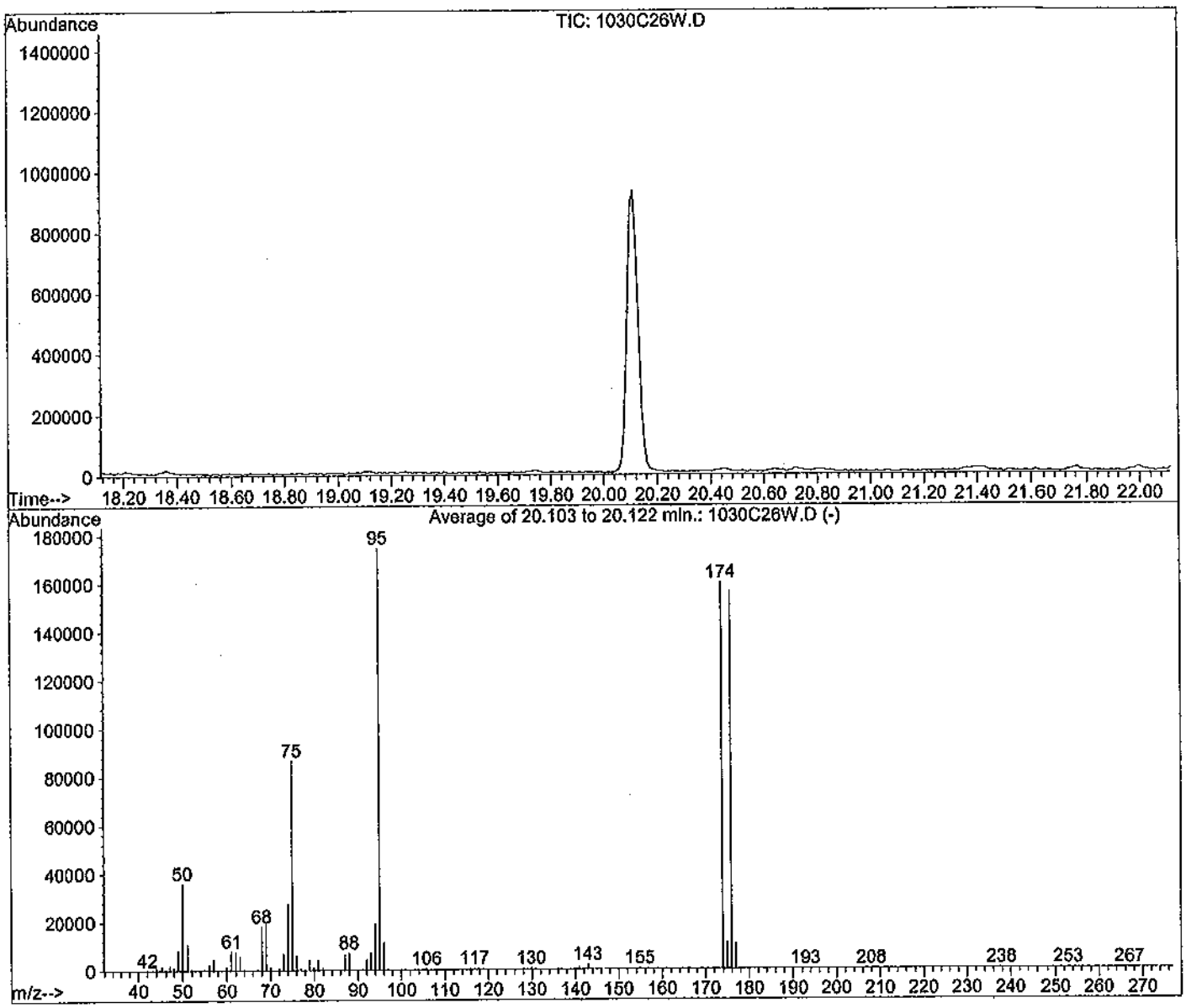
Target Mass	Rel. to Mass	Lower Limit%	Upper Limit%	Rel. Abn%	Raw Abn	Result Pass/Fail
50	95	15	40	17.7	88217	PASS
75	95	30	60	45.1	224883	PASS
95	95	100	100	100.0	499051	PASS
96	95	5	9	6.5	32634	PASS
173	174	0.00	2	0.0	0	PASS
174	95	50	100	92.8	463189	PASS
175	174	5	9	7.2	33219	PASS
176	174	95	101	97.1	449771	PASS
177	176	5	9	6.4	28567	PASS

1030C13W.D CALLW.M Tue Dec 06 16:03:25 2011

Data File : M:\CHICO\DATA\C111030\1030C26W.D
Acq On : 31 Oct 11 7:21
Sample : 20ug/mL BFB STD 10-19-11
Misc : Water 2ul

Vial: 1
Operator: STC
Inst : Chico
Multiplr: 1.00

Method : M:\CHICO\DATA\C111030\CALLW.M (RTE Integrator)
Title : METHOD 8260



Spectrum Information: Average of 20.103 to 20.122 min.

Target Mass	Rel. to Mass	Lower Limit%	Upper Limit%	Rel. Abn%	Raw Abn	Result Pass/Fail
50	95	15	40	20.6	35941	PASS
75	95	30	60	49.7	86725	PASS
95	95	100	100	100.0	174635	PASS
96	95	5	9	6.6	11458	PASS
173	174	0.00	2	0.0	0	PASS
174	95	50	100	91.6	159913	PASS
175	174	5	9	6.9	11031	PASS
176	174	95	101	97.7	156203	PASS
177	176	5	9	6.7	10398	PASS

8-15-11
RS

Volatile Standard Curve Preparation for 5mL Purge (8260 soil)-THOR

	Expiration Date:		08/16/11							
		5µg/mL Vol Std #9	5µg/mL Surr	50µg/mL Vol Std #7	50µg/mL Vol Std #8	50µg/mL Surr	5µg/mL Vol Std #10	50µg/mL Vol Std #1	5µg/mL Vol Std #2	50µg/mL Vol Std #12
Date	Conc.	08-10-11R	08-10-11V	08-10-11N	08-10-11P	08-10-11U	08-10-11S	08-10-11Q	08-10-11Q	08-10-11T
Code	µg/L	Exp:08-17-11	Exp:08-17-11	Exp:08-17-11	Exp:08-17-11	Exp:08-17-11	Exp:08-17-11	Exp:08-17-11	Exp:08-17-11	Exp:08-17-11
08-15-11A	2	2	2	n/a	n/a	n/a	2	n/a	2	n/a
08-15-11B	5	5	5	n/a	n/a	n/a	5	n/a	5	n/a
08-15-11C	10	10	10	n/a	n/a	n/a	10	n/a	10	n/a
08-15-11D	20	20	20	n/a	n/a	n/a	20	n/a	20	n/a
08-15-11E	50	n/a	n/a	5	5	5	n/a	5	n/a	5
08-15-11F	100	n/a	n/a	10	10	10	n/a	10	n/a	10
08-15-11G	200	n/a	n/a	20	20	20	n/a	20	n/a	20

250µg/mL TBA	Final Vol
08-10-11W	w/P&T H2O
Exp:08-17-11	mL
1	5
2	5
3	5
4	5
5	5
6	5
7	5

8-16-11
RS

Volatile Standard Curve Preparation for 5mL Purge (8260 soil)-THOR

	Expiration Date:		08/17/11							
		5µg/mL Vol Std #9	5µg/mL Surr	50µg/mL Vol Std #7	50µg/mL Vol Std #8	50µg/mL Surr	5µg/mL Vol Std #10	50µg/mL Vol Std #1	5µg/mL Vol Std #2	50µg/mL Vol Std #12
Date	Conc.	08-10-11R	08-10-11V	08-10-11N	08-10-11P	08-10-11U	08-10-11S	08-10-11Q	08-10-11Q	08-10-11T
Code	µg/L	Exp:08-17-11	Exp:08-17-11	Exp:08-17-11	Exp:08-17-11	Exp:08-17-11	Exp:08-17-11	Exp:08-17-11	Exp:08-17-11	Exp:08-17-11
08-16-11A	2	2	2	n/a	n/a	n/a	2	n/a	2	n/a
08-16-11B	5	5	5	n/a	n/a	n/a	5	n/a	5	n/a
08-16-11C	10	10	10	n/a	n/a	n/a	10	n/a	10	n/a
08-16-11D	20	20	20	n/a	n/a	n/a	20	n/a	20	n/a
08-16-11E	50	n/a	n/a	5	5	5	n/a	5	n/a	5
08-16-11F	100	n/a	n/a	10	10	10	n/a	10	n/a	10
08-16-11G	200	n/a	n/a	20	20	20	n/a	20	n/a	20

250µg/mL TBA	Final Vol
08-10-11W	w/P&T H2O
Exp:08-17-11	mL
1	5
2	5
3	5
4	5
5	5
6	5
7	5

8-19-11
A -
RS

Method 8260 Gases, 2,000 mg/L, 2 X 0.6 ml

120016-03

Lot #	Storage	Expiry
170302	≤-10 Degrees C	3/11/14

Solvt: P/T Methanol

Method 8260 Gases
Lot #: 170302 - 28666
Rec: 4/20/11 MFR exp. 03/11/14

8-19-11
B -
RS

Method 8260B Surrogate Solution, 2,000 mg/L, 1 ml

120002-01

Lot #	Storage	Expiry
164585	-10 Degrees C	10/12/13

Solvt: P/T Methanol

Method 8260B Surrogate
Lot #: 164585 - 28720
Rec: 4/20/11 MFR exp. 10/12/13

8-19-11
C -
RS

VOC Mix 4-3, 2,000 mg/L, 1 ml

120166-01

Lot #	Storage	Expiry
171714	≤-6 Degrees C	4/11/13

Solvt: P/T Methanol

VOC Mix 4-3, 2000mg/L
Lot #: 171714 - 29243
Rec: 8/5/11 MFR exp. 04/11/13

GC/MS STANDARD PREPARATION BOOK # 58 PAGE # 4

		08-17-11V	Exp:	08/24/11			
		5ug/ml Vol Work Std #9					
		SOURCES	Lot	APPL Code	APPL Exp Date	ul	
		50ug/ml Vol Work Std #7		08-17-11R	09/02/11	200	
		50ug/ml Vol Work Std #8		08-17-11T	09/02/11	200	
		J&T Brand		08/12/12	06/08/12	1600	
		08-17-11W	Exp:	08/24/11			
		5ug/ml Vol Work Std #10					
		SOURCES	Lot	APPL Code	APPL Exp Date	ul	
		50ug/ml Vol Work Std #1		08-17-11S	09/02/11	200	
		J&T Brand		08/12/12	06/08/12	1800	
		08-17-11X	Exp:	08/24/11			
		5ug/ml Vol Work Std #12					
		SOURCES	Lot	APPL Code	APPL Exp Date	ul	
		50ug/ml Vol Work Std #2		08-17-11U	09/02/11	200	
		J&T Brand		08/12/12	06/08/12	1800	
08-17-11Y							
50ug/ml 8260 Surrogate			Conc.		Date	Exp.	
Exp:08/24/11			ug/ml	Lot #	Code	Date	ul
02SI	120002-01	8260B Surr Solution	2000	164585-28720	08-17-11B	09/14/11	100
J&T Brand		Purge & Trap MeOH		K07E34-00543	08/12/12	10/14/11	1900
08-17-11Z			Exp:	08/24/11			
5.0ug/ml 8260 Surrogate			Lot	APPL Code	APPL Exp Date	ul	
		50ug/ml 8260 Surrogate		08-17-11Y	09/02/11	200	
J&T Brand		Purge & Trap MeOH		08/05/12	06/08/12	1800	
08-17-11AA							
250ug/ml TBA/IBA/Acetonitrile/Cyclohexanone/Acrolein/2-P						APPL	
Exp:08/24/11			Conc.		Date	Exp.	
Supplier	ID #		ug/ml	Lot #	Code	Date	ul
02SI	120166-01	Volatile Mix 4-3	2000	171714-29243	08-17-11C	12/17/11	500
02SI	020229-09	Acrolein	10000	175935-29032	08-04-11J	08/22/11	100
J&T Brand		Purge & Trap MeOH		K07E14-00543	08/12/12	10/14/11	1400

8-17-11 RS

8-17-11 RS

8-17-11 RS

8-17-11 RS

CHICO							
08-17-11AB							
250ug/ml 8260 Internal Standard - Chico			Conc.		Date	Exp.	
Supplier	ID #		ug/ml	Lot #	Code	Date	ul
02SI	120302-03	Internal Standard Mix	2000	166255-27947	08-09-11A	10/23/11	500
02SI	020132-02	Fluorobenzene Standard	2000	169170-28263	08-09-11B	10/23/11	500
J&T Baker		Purge & Trap MeOH		K07E34-00543	08/12/11	11/14/12	5000
08-17-11AC							
250ug/ml 8260 Surrogate - Chico			Conc.		Date	Exp.	
Supplier	ID #		ug/ml	Lot #	Code	Date	ul
02SI	120002-01	Surrogate Standard	2000	164585-28727	08-09-11C	10/23/11	500
J&T Baker		Purge & Trap MeOH		K07E34-00543	08/12/11	11/14/12	5000

8-17-11 RS

8-18-11 RS

08-17-11AD							
50ug/ml 8260B Surrogate- Nsc			Conc.		Date	Exp.	
Supplier	ID #		ug/ml	Lot #	Code	Date	ul
02SI	8260B Surr	Surrogate Standards	2000	164585-28720	08-17-11B	12/13/11	500
J.T Baker		Purge & Trap MeOH		K07E34-00543	08/12/11	10/10/11	19500

Volatile Standard Curve Preparation for 10mL Purge (8260 water)-THOR

	Expiration Date:		09/22/11								
		5ug/mL Vol Std #9	5ug/mL Surr	50ug/mL Vol Std #7	50ug/mL Vol Std #8	50ug/mL Surr	5ug/mL Vol Std #10	50ug/mL Vol Std #6	5ug/mL Vol Std #12	250ug/mL TAPD
Date	Conc.	09-19-11I	09-19-11M	09-19-11O	09-19-11F	09-19-11L	09-19-11J	09-19-11E	09-19-11G	09-19-11K
Code	µg/L	Exp:09-26-11	Exp:09-26-11	Exp:09-26-11	Exp:09-26-11	Exp:09-26-11	Exp:09-26-11	Exp:09-26-11	Exp:09-26-11	Exp:09-26-11
09-21-11B	0.3	3	6	n/a	n/a	n/a	n/a	n/a	n/a	3
09-21-11C	0.5	5	10	n/a	n/a	n/a	5	n/a	n/a	5
09-21-11D	1	10	20	n/a	n/a	n/a	10	n/a	n/a	10
09-21-11E	2	20	40	n/a	n/a	n/a	20	n/a	n/a	20
09-21-11F	5	n/a	n/a	5	5	10	n/a	5	5	n/a
09-21-11G	10	n/a	n/a	10	10	25	n/a	10	10	n/a
09-21-11H	20	n/a	n/a	20	20	40	n/a	20	20	n/a
09-21-11I	40	n/a	n/a	40	40	80	n/a	40	40	n/a
09-21-11J	100	n/a	n/a	100	100	n/a	n/a	100	100	n/a

250ug/mL TAPD 09-19-11K Exp:09-26-11	Final Vol w/P&T H2O mL
3	50
5	50
10	50
15	50
20	50
25	50
30	50
35	50
40	50

Volatile Standard Curve Preparation for 10mL Purge (524 water)-NEO

	Expiration Date:		09/22/11				
		5ug/mL Vol Std #9	50ug/mL Vol Std #7	50ug/mL Vol Std #8	250ug/mL TAPD	Final Vol
Date	Conc.	09-19-11I	09-19-11O	09-19-11F	09-19-11N	w/P&T H2O
Code	µg/L	Exp:09-26-11	Exp:09-26-11	Exp:09-26-11	Exp:09-26-11	mL
09-21-11K	0.2	2	n/a	n/a	2	50
09-21-11L	0.5	5	n/a	n/a	5	50
09-21-11M	1	10	n/a	n/a	10	50
09-21-11N	2	20	n/a	n/a	15	50
09-21-11O	5	n/a	5	5	20	50
09-21-11P	10	n/a	10	10	25	50
09-21-11Q	40	n/a	40	40	35	50
09-21-11R	100	n/a	100	100	40	50

AY44799W01

26

VOLATILES 38083 - 062910

65441
AY44799 W01
VOA_Frig

Part #: 38083 — Laboratory Use Only - See MSDS
Lot #: 062910 — Exp: 062913 — Storage 0 °C
CWA Volatiles in Non-Potable Water
Varied in methanol — 2 mL
ABSOLUTE STANDARDS, INC. • 800-368-1131

Sweetpea

09-23-11A

250ug/ml 8260 Internal Standard - Sweetpea			Conc.		Date	Exp.	
Supplier	ID #		ug/ml	Lot #	Code	Date	uL
02SI	120302-03	Internal Standard Mix	2000	166255-28349	09-14-11A	10/10/11	500
	020132-02	Fluorobenzene Standard	2000	169170-28739	09-14-11B	10/10/11	500
J.T.Baker		Purge & Trap MeOH		K07E34-00547	09/12/11	10/14/12	3000

Method 8260 Internal Standard Solution, 2,000 mg/L, 1 ml
120302-03
Lot # 166255 — Storage ≤ 10 Degrees C — Expiry 10/10/11
Solv: P/T Methanol
Made in the USA

Method 8260 Internal Standard
Lot #: 166255 - 28350
Rec: 2/17/11 MFR exp. 11/18/12

Fluorobenzene Solution, 2,000 mg/L, 1 ml
020132-02
Lot # 169170 — Storage ≤ 6 Degrees C — Expiry 2/13/14
Solv: P/T Methanol
Made in the USA

Fluorobenzene
Lot #: 169170 - 28738
Rec: 4/20/11 MFR exp. 02/13/14

CHICO					
09-23-11D					
250ug/ml 8260 Internal Standard - Chico			Conc.		Date
Supplier	ID #		ug/ml	Lot #	Code
02SI	120302-03	Internal Standard Mix	2000	166255-28350	09-23-11B
02SI	020132-02	Fluorobenzene Standard	2000	169170-28738	09-23-11C
J&T Baker		Purge & Trap MeOH		K07B34-00547	09/12/11

9.23.11 RS.

Volatile Standard Curve Preparation for 10mL Purge (8260 water)-SWEETPEA

	Expiration Date:		09/24/11							
		5µg/mL Vol Std #9	5µg/mL Surr	50µg/mL Vol Std #7	50µg/mL Vol Std #8	50µg/mL Surr	5µg/mL Vol Std #10	50µg/mL Vol Std #1	50µg/mL Vol Std #2
Date	Conc.	09-19-11I	09-19-11M	09-19-11D	09-19-11F	09-19-11L	09-19-11J	09-19-11E	09-19-11G
Code	µg/L	Exp:09-26-11	Exp:09-26-11	Exp:09-26-11	Exp:09-26-11	Exp:09-26-11	Exp:09-26-11	Exp:09-26-11	Exp:09-26-11
09-23-11E	0.3	3	6	n/a	n/a	n/a	3	n/a	n/a
09-23-11F	0.5	5	10	n/a	n/a	n/a	5	n/a	n/a
09-23-11G	1	10	20	n/a	n/a	n/a	10	n/a	n/a
09-23-11H	2	20	40	n/a	n/a	n/a	20	n/a	n/a
09-23-11I	5	n/a	n/a	5	5	10	n/a	5	5
09-23-11J	10	n/a	n/a	10	10	25	n/a	10	10
09-23-11K	20	n/a	n/a	20	20	40	n/a	20	20
09-23-11L	40	n/a	n/a	40	40	80	n/a	40	40
09-23-11M	100	n/a	n/a	100	100	n/a	n/a	100	100

9.23.11 RS.

9.24.11 RS.

Volatile Standard Curve Preparation for 10mL Purge (8260 water)-CHICO

	Expiration Date:		09/24/11							
		5µg/mL Vol Std #9	5µg/mL Surr	50µg/mL Vol Std #7	50µg/mL Vol Std #8	50µg/mL Surr	5µg/mL Vol Std #10	50µg/mL Vol Std #1	50µg/mL Vol Std #2
Date	Conc.	09-19-11I	09-19-11M	09-19-11D	09-19-11F	09-19-11L	09-19-11J	09-19-11E	09-19-11G
Code	µg/L	Exp:09-26-11	Exp:09-26-11	Exp:09-26-11	Exp:09-26-11	Exp:09-26-11	Exp:09-26-11	Exp:09-26-11	Exp:09-26-11
09-23-11N	0.3	3	6	n/a	n/a	n/a	3	n/a	n/a
09-23-11O	0.5	5	10	n/a	n/a	n/a	5	n/a	n/a
09-23-11P	1	10	20	n/a	n/a	n/a	10	n/a	n/a
09-23-11Q	2	20	40	n/a	n/a	n/a	20	n/a	n/a
09-23-11R	5	n/a	n/a	5	5	10	n/a	5	5
09-23-11S	10	n/a	n/a	10	10	25	n/a	10	10
09-23-11T	20	n/a	n/a	20	20	40	n/a	20	20
09-23-11U	40	n/a	n/a	40	40	80	n/a	40	40
09-23-11V	100	n/a	n/a	100	100	n/a	n/a	100	100

9.23.11 RS.

9.24.11 RS.

9.24.11 A- RS.

Method 8260 Gases, 2,000 mg/L, 2 X 0.5 ml

120016-05

Lot #	Storage	Expiry
170302	≤-10 Degrees C	3/11/14

Solvt P/T Methanol

Method 8260 Gases

Lot #: 170302 - 28677

Rec: 4/20/11 MFR exp. 03/11/14

RS.

9.24.11 B- RS.

n-Hexane Solution, 1,000 mg/L, 1 ml

020610-02

Lot #	Storage	Expiry
163378	≤-10 Degrees	8/29/15

Solvt P/T Methanol

n-Hexane Solution

Lot #: 163378 - 29230

Rec: 8/5/11 MFR exp. 08/29/15

RS.

9.24.11 RS.

9.24.11 RS.

068

GC/MS STANDARD PREPARATION BOOK # _____ PAGE # _____

Volatile Standard Curve Preparation for 10mL Purge (5250 water)-CHICO

	Expiration Date:		09/30/11							
		5µg/mL Vol Std #9	5µg/mL Surr	50µg/mL Vol Std #7	50µg/mL Vol Std #8	50µg/mL Surr	5µg/mL Vol Std #10	50µg/mL Vol Std #11	50µg/mL Vol Std #12	250µg/mL TAPD
Date	Conc.	09-29-11D	09-29-11N	09-29-11F	09-29-11C	09-29-11M	09-29-11K	09-29-11G	09-29-11I	09-29-11O
Code	µg/L	Exp:10-05-11	Exp:10-05-11	Exp:10-05-11	Exp:10-05-11	Exp:10-05-11	Exp:10-05-11	Exp:10-05-11	Exp:10-05-11	Exp:10-05-11
09-29-11N	0.3	3	6	n/a	n/a	n/a	3	n/a	n/a	3
09-29-11O	0.5	5	10	n/a	n/a	n/a	5	n/a	n/a	5
09-29-11P	1	10	20	n/a	n/a	n/a	10	n/a	n/a	10
09-29-11Q	2	20	40	n/a	n/a	n/a	20	n/a	n/a	20
09-29-11R	5	n/a	n/a	5	5	10	n/a	5	5	5
09-29-11S	10	n/a	n/a	10	10	25	n/a	10	10	10
09-29-11T	20	n/a	n/a	20	20	40	n/a	20	20	20
09-29-11U	40	n/a	n/a	40	40	60	n/a	40	40	40
09-29-11V	100	n/a	n/a	100	100	n/a	100	n/a	100	100

9-29-11 RS.

10-02-11 RS

10-02-11 RS

10-02-11 RS

10-02-11 RS

10-02-11 RS

4-Bromofluorobenzene
Solution, 2,000 mg/L, 1 ml
Lot # 020135-03
Expiry
4-Bromofluorobenzene
Lot #: 176675-29375
Rec: 8/9/11 MFR exp. 08/02/14

9-30-11 A- RS.

9-30-11 RS.

09-30-11A							
20ug/ml BFB STD			Conc.		Date	EXP:	
EXP: 10-30-11			ug/ml	Lot#	CODE	Date:	
O2SI	020135-03	4-Bromofluorobenzene	2500	176675-29375	10-01-11A	12/11/14	
J&T Baker		Purge & Trap MeOH		K14E06-00551	10/01/11	09/29/12	

Volatile Standard Curve Preparation for 10mL Purge (524 water)-NEO

	Expiration Date:		10/02/11				
		5µg/mL Vol Std #9	50µg/mL Vol Std #7	50µg/mL Vol Std #8	250µg/mL TAPD	Final Vol
Date	Conc.	09-29-11D	09-29-11F	09-29-11C	09-29-11O	wP&T H2O
Code	µg/L	Exp:10-05-11	Exp:10-05-11	Exp:10-05-11	Exp:10-05-11	mL
10-01-11A	0.2	2	n/a	n/a	2	50
10-01-11B	0.5	5	n/a	n/a	5	50
10-01-11C	1	10	n/a	n/a	10	50
10-01-11D	2	20	n/a	n/a	15	50
10-01-11E	5	n/a	5	5	20	50
10-01-11F	10	n/a	10	10	25	50
10-01-11G	40	n/a	40	40	35	50
10-01-11H	100	n/a	100	100	40	50

10-01-11 RS.

10-02-11 RS

Volatile Standard Curve Preparation for 10mL Purge (8260 water)-MAX

Expiration Date: 10/28/2011

Date	Conc.	5µg/mL Vol Std #9	5µg/mL Surr	50µg/mL Vol Std #7	50µg/mL Vol Std #8	50µg/mL Surr	5µg/mL Vol Std #10	50µg/mL Vol Std #1	5µg/mL Vol Std #2	5µg/mL Vol Std #12
Code	µg/L	10-26-11J Exp:11-01-11	10-26-11N Exp:11-01-11	10-26-11F Exp:11-01-11	10-26-11H Exp:11-01-11	10-26-11M Exp:11-01-11	10-26-11K Exp:11-01-11	10-26-11G Exp:11-01-11	10-26-11I Exp:11-01-11	10-26-11L Exp:11-01-11
10-27-11J	0.5	3	6	n/a	n/a	n/a	3	n/a	n/a	3
10-27-11K	0.5	5	10	n/a	n/a	n/a	5	n/a	n/a	5
10-27-11L	1	10	20	n/a	n/a	n/a	10	n/a	n/a	10
10-27-11M	2	20	40	n/a	n/a	n/a	20	n/a	n/a	20
10-27-11N	5	n/a	n/a	5	5	10	n/a	5	5	n/a
10-27-11O	10	n/a	n/a	10	10	25	n/a	10	10	n/a
10-27-11P	20	n/a	n/a	20	20	40	n/a	20	20	n/a
10-27-11Q	40	n/a	n/a	40	40	80	n/a	40	40	n/a
10-27-11R	100	n/a	n/a	100	100	n/a	n/a	100	100	n/a

10-27-11
RS

250µg/mL TAPD 10-26-11O Exp:11-01-11	Final Vol w/P&T H2O mL
3	50
5	50
10	50
15	50
20	50
25	50
30	50
35	50
40	50

Volatile Standard Curve Preparation for 5mL Purge (8260 soil)-NEO

Expiration Date: 10/28/2011

Date	Conc.	5µg/mL Vol Std #9	5µg/mL Surr	50µg/mL Vol Std #7	50µg/mL Vol Std #8	50µg/mL Surr	5µg/mL Vol Std #10	50µg/mL Vol Std #1	5µg/mL Vol Std #2	50µg/mL Vol Std #12
Code	µg/L	10-26-11J Exp:11-01-11	10-26-11N Exp:11-01-11	10-26-11F Exp:11-01-11	10-26-11H Exp:11-01-11	10-26-11M Exp:11-01-11	10-26-11K Exp:11-01-11	10-26-11G Exp:11-01-11	10-26-11I Exp:11-01-11	10-26-11L Exp:11-01-11
10-27-11S	2	2	2	n/a	n/a	n/a	2	n/a	2	n/a
10-27-11T	5	5	5	n/a	n/a	n/a	5	n/a	5	n/a
10-27-11U	10	10	10	n/a	n/a	n/a	10	n/a	10	n/a
10-27-11V	20	20	20	n/a	n/a	n/a	20	n/a	20	n/a
10-27-11W	50	n/a	n/a	5	5	5	n/a	5	n/a	5
10-27-11X	100	n/a	n/a	10	10	10	n/a	10	n/a	10
10-27-11Y	200	n/a	n/a	20	20	20	n/a	20	n/a	20

10-27-11
RS

250µg/mL TBA 10-26-11O Exp:11-01-11	Final Vol w/P&T H2O mL
1	5
2	5
3	5
4	5
5	5
6	5
7	5

Volatile Standard Curve Preparation for 5mL Purge (8260 soil)-THOR

Expiration Date: 10/28/2011

Date	Conc.	5µg/mL Vol Std #9	5µg/mL Surr	50µg/mL Vol Std #7	50µg/mL Vol Std #8	50µg/mL Surr	5µg/mL Vol Std #10	50µg/mL Vol Std #1	5µg/mL Vol Std #2	50µg/mL Vol Std #12
Code	µg/L	10-26-11J Exp:11-01-11	10-26-11N Exp:11-01-11	10-26-11F Exp:11-01-11	10-26-11H Exp:11-01-11	10-26-11M Exp:11-01-11	10-26-11K Exp:11-01-11	10-26-11G Exp:11-01-11	10-26-11I Exp:11-01-11	10-26-11L Exp:11-01-11
10-27-11Z	2	2	2	n/a	n/a	n/a	2	n/a	2	n/a
10-27-11AA	5	5	5	n/a	n/a	n/a	5	n/a	5	n/a
10-27-11AB	10	10	10	n/a	n/a	n/a	10	n/a	10	n/a
10-27-11AC	20	20	20	n/a	n/a	n/a	20	n/a	20	n/a
10-27-11AD	50	n/a	n/a	5	5	5	n/a	5	n/a	5
10-27-11AE	100	n/a	n/a	10	10	10	n/a	10	n/a	10
10-27-11AF	200	n/a	n/a	20	20	20	n/a	20	n/a	20

6/27/11
RS.

250µg/mL TBA 10-26-11O Exp:11-01-11	Final Vol w/P&T H2O mL
1	5
2	5
3	5
4	5
5	5
6	5
7	5

Method 8260 Gases, 2,000 mg/L, 2 X 0.6 ml

I20016-03

Lot #	Storage	Expiry
169238	≤ -10 Degrees C	2/19/14

Solv: P/T Methanol

Method 8260 Gases

Lot #: 169238 - 28682

Rec: 4/20/11 MFR exp. 02/19/14

10-28-11 A-
RS.

RS.

10-28-11 RS

B -

Hexachloroethane Solution, 1000 mg/L, 1 ml

020049-02

Lot #	Storage	Expiry
164816	≤ -10 Degrees C	10/14/12

Solv: P/T Methanol

Hexachloroethane

Lot #: 164816 - 28687

Rec: 4/20/11 MFR exp. 10/14/12

RS

10-28-11 RS.

B

10-28-11 RS

C -

Benzyl Chloride Solution, 1000 mg/L, 1 ml

020228-02

Lot #	Storage	Expiry
163373	≤ -10 Degrees	8/29/11

Solv: P/T Methanol

Benzyl Chloride

Lot #: 163373 - 29166

Rec: 8/5/11 MFR exp. 08/29/12

RS

10-28-11 RS.

C

10-28-11 RS

D -

Volatile Mix, 20-29, 2,000 mg/L, 1 ml

121039-02

Lot #	Storage	Expiry
163374	≤ -10 Degrees	8/29/12

Solv: P/T Methanol

Volatile Mix, 20-29

Lot #: 163374 - 28300

Rec: 2/17/11 MFR exp. 08/29/12

RS

10-28-11 RS.

D

10-28-11 RS

E -

Method 8260 VOC Liquids, 54 Compounds, 2,000 mg/L, 1 ml

120023-03

Lot #	Storage	Expiry
164454	≤ -10 Degrees C	10/4/12

Solv: P/T Methanol

8260 VOC Liquids, 54 Comp.

Lot #: 164454 - 27872

Rec: 12/15/10 MFR exp. 10/04/12

RS

10-28-11 RS.

E

10-28-11 RS

F -

Vinyl Acetate Solution, 2,000 mg/L, 1ml

Lot #	Storage	Expiry
178902	≤ -10 Degrees C	12/15/11

Solv: P/T Methanol

Vinyl Acetate

Lot #: 178902 - 29552

Rec: 9/22/11 MFR exp. 12/15/11

RS

10-28-11 G -
RS.

Heptane Solution, 1000 mg/L, 1 ml
R&D only, not human consump. Made in the USA
820546-82
Lot # 169174 | Storage ≤-10 Degrees C | Expiry 2/18/14
Solv: P/T Methanol

RS

Heptane Solution
Lot #: 169174 - 29248
Rec: 8/5/11 MFR exp. 02/18/14

10-28-11 H -
RS.

8260B Surrogate Solution, 2,000 mg/L, 5 x 1 ml
R&D only, not human consump. Made in the USA
120002-01-5PAK
Lot # 173249 | Storage ≤-10 Degrees C | Expiry 5/17/13
Solv: P/T Methanol

RS.

8260B Surrogate Solution
Lot #: 173249 - 28847
Rec: 5/25/11 MFR exp. 05/17/13

10-28-11 I -
RS.

VOC Mix 4-3, 2,000 mg/L, 1 ml
Made in the USA
120166-01
Lot # 178651 | Storage ≤-10 Degrees C | Expiry 9/11/13
Solv: P/T Methanol

RS.

VOC Mix 4-3, 2000mg/L
Lot #: 178651 - 29510
Rec: 9/20/11 MFR exp. 09/11/13

10-28-11 J -
RS.

Method 8260 Gases (Second Source), 2,000 mg/L, 2 X 0.6 ml
R&D only, not human consump. Made in the USA
120016-03-SS
Lot # 168038 | Storage ≤-10 Degrees C | Expiry 1/21/14
Solv: P/T Methanol

RS.

8260 Gases (SS)
Lot #: 168038 - 28743
Rec: 4/20/11 MFR exp. 01/21/14

10-28-11K							
50ug/ml Vol Work Std #7							
Exp:11/04/11							
			Conc.		Date	Exp.	
Supplier	ID #	ID	ug/ml	Lot #	Code	Date	ul
02SI	120016-01	Gas Mix	2000	169238-28582	10-28-11A	11/30/2011	100
02SI	020049-02	HEXACHLOROETHANE	1000	164816-28687	10-28-11B	12/14/2011	200
02SI	020228-02	Benzyl Chloride	1000	163173-29166	10-28-11C	12/14/2011	200
J&T Brand		Purge & Trap MeOH		K14E06-00556	10/27/2011	6/8/2012	3500
10-28-11L							
50ug/ml Vol Work Std #1							
Exp:11/04/11							
Supplier	ID #	ID	ug/ml	Lot #	Code	Date	ul
02SI	020145-02-02	2-CEVE	2000	160092-26537	10-06-11B	12/7/2011	50
J&T Brand		Purge & Trap MeOH		K14E06-00556	10/27/2011	6/8/2012	1950
10-28-11M							
50ug/ml Vol Work Std #8							
Exp:11/04/11							
			Conc.		Date	Exp.	
Supplier	ID #	ID	ug/ml	Lot #	Code	Date	ul
02SI	122039-02	Volatile Mix, 20-29	2000	163374-28100	10-28-11D	2/14/2012	100
02SI	120023-03	VOC'S-54 COMP	2000	164454-27872	10-28-11E	2/14/2012	100
02SI	020232-02	Vinyl Acetate	2000	178902-29552	10-28-11F	11/15/2011	100
02SI	020620-02	n-Hexane	1000	163378-27889	10-26-11B	11/14/2011	200
02SI	020546-02	Heptane	1000	169174-29248	10-28-11G	11/14/2011	200
J&T Brand		Purge & Trap MeOH		K14E06-00556	10/27/2011	6/8/2012	3300
10-28-11N							
50ug/ml Vol Work Std #2							
Exp:11/04/11							
Supplier	ID #	ID	ug/ml				
02SI	121020-05	HSL'S-Ketone Solution	2000	169173-28307	10-12-11E	11/14/2011	100
J&T Brand		Purge & Trap MeOH		K14E06-00556	10/27/2011	10/14/2012	3900
		10-28-11O	Exp:	11/4/2011			
		5ug/ml Vol Work Std #3					
		SOURCES	Lot	APPL Code	APPL Exp Date	ul	
		50ug/ml Vol Work Std #7		10-28-11K	10/31/2011	200	
		50ug/ml Vol Work Std #8		10-28-11M	10/31/2011	200	
		J&T Brand		10/6/2011	6/8/2012	1600	
		10-28-11P	Exp:	11/4/2011			
		5ug/ml Vol Work Std #10					
		SOURCES	Lot	APPL Code	APPL Exp Date	ul	
		50ug/ml Vol Work Std #1		10-28-11L	10/31/2011	200	
		J&T Brand		10/27/2011	6/8/2012	1800	
		10-28-11Q	Exp:	11/4/2011			
		5ug/ml Vol Work Std #12					
		SOURCES	Lot	APPL Code	APPL Exp Date	ul	
		50ug/ml Vol Work Std #2		10-28-11N	10/31/2011	200	
		J&T Brand		10/27/2011	6/8/2012	1800	
10-28-11R							
50ug/ml 8260 Surrogate			Conc.		Date	Exp.	
Exp:11/04/11			ug/ml	Lot #	Code	Date	ul
02SI	120002-01	8260B Surr Solution	2000	173249-28847	10-28-11H	11/14/2011	100
J&T Brand		Purge & Trap MeOH		K14E06-00556	10/27/2011	6/8/2012	3900
10-28-11S			Exp:	11/4/2011			
5.0ug/ml 8260 Surrogate			Lot	APPL Code	APPL Exp Date	ul	
		50ug/ml 8260 Surrogate		10-28-11R	10/31/2011	200	
J&T Brand		Purge & Trap MeOH		K14E06-00556	10/27/2011	1800	
10-28-11T							
250ug/ml TBA/IPA/Acetonitrile/Cyclohexanone/Acrolein/2-P						APPL	
Exp:11/04/11			Conc.		Date	Exp.	
Supplier	ID #		ug/ml	Lot #	Code	Date	ul
02SI	120166-01	Volatile Mix 4-3	2000	178691-29510	10-28-11I	12/17/2011	500
02SI	020228-09	Acrolein	10000	179941-29561	10-19-11H	11/21/2011	100
J&T Brand		Purge & Trap MeOH		K14E06-00556	10/27/2011	6/8/2012	3400

10/28/11 RS.

10/28/11 RS.

10/28/11 RS.

10-28-11U							
50ug/ml VOC Std#5							
Exp:11/04/11							
			Conc.		Date	Exp.	
Supplier	ID #	ID	ug/ml	Lot #	Code	Date	ul
O2SI	120016-03-SS	8260 Gases(SS)	2000	168038-28741	10-28-11J	11/30/2011	50
O2SI	020145-02-02-S	2-CEVE	2000	162530-27271	10-19-11J	11/3/2011	50
J&T Brand		Purge & Trap MeOH		KL4E06-00556	10/27/2011	6/8/2012	1900
10-28-11V							
50ug/ml VOC Std#6							
Exp:11/04/11							
	ID #	ID	ug/ml	Lot #	Code	Date	ul
O2SI	120021-03-SS	VOC'S 54 COMP.	2000	161271-27771	09-12-11P	11/14/2011	50
O2SI	120296-01	Custom 8260 Solution	2000	166038-27763	09-12-11Q	11/14/2011	50
O2SI	020232-02-SS	Vinyl Acetate(SS)	2000	176774-29257	09-12-11R	11/30/2011	50
O2SI	020620-02-SS	n-HEXANE	1000	179199-29615	10-12-11F	12/14/2011	100
O2SI	020049-02-SS	HEXACHLOROETHANE	1000	154535-25913	09-13-11B	12/29/2011	100
O2SI	020546-02-SS	Heptane(SS)	1000	142276-23593	09-13-11C	12/19/2011	100
J&T Brand		Purge & Trap MeOH		K14E06-00556	10/27/2011	6/8/2012	1550
10-28-11W							
250ug/ml TBA/IBA/Acetonitrile/Cyclohexanone/Acrolein/2-P						APPL	
Exp:11/04/11						Exp.	
			Conc.		Date		
Supplier	ID #	ID	ug/ml	Lot #	Code	Date	uL
O2SI	120166-01-SS	VOC Mix 4-3 (SS)	2000	152531-25468	10-02-11G	11/3/2011	250
O2SI	020229-09-SS	Acrolein SOLUTION (SS)	10000	178607-29549	10-02-11H	11/21/2011	50
J&T Brand		Purge & Trap MeOH		K14E06-00556	10/27/2011	6/8/2012	1700

10-28-11X							
50ug/ml Vol Work Std #7							
Exp:11/04/11							
			Conc.		Date	Exp.	
Supplier	ID #	ID	ug/ml	Lot #	Code	Date	ul
O2SI	120016-03	Gas Mix	2000	169238-28682	10-28-11A	11/30/2011	100
O2SI	020049-02	HEXACHLOROETHANE	1000	164816-28687	10-28-11B	12/14/2011	200
O2SI	020228-02	Benzyl Chloride	1000	163371-29166	10-28-11C	12/14/2011	200
J&T Brand		Purge & Trap MeOH		K14E06-00556	10/27/2011	6/8/2012	3500
10-28-11Y							
50ug/ml Vol Work Std #1							
Exp:11/04/11							
Supplier	Id #	ID	ug/ml	Lot #	Code	Date	ul
O2SI	020145-02-02	2-CEVE	2000	160092-26617	10-06-11B	12/7/2011	50
J&T Brand		Purge & Trap MeOH		K14E06-00556	10/27/2011	6/8/2012	1950
10-28-11Z							
50ug/ml Vol Work Std #8							
Exp:11/04/11							
			Conc.		Date	Exp.	
Supplier	ID #	ID	ug/ml	Lot #	Code	Date	ul
O2SI	122039-02	Volatile Mix, 20-29	2000	163374-28300	10-28-11D	2/14/2012	100
O2SI	120023-01	VOC'S-54 COMP	2000	164454-27872	10-28-11E	2/14/2012	100
O2SI	020232-02	Vinyl Acetate	2000	178902-29552	10-28-11F	11/15/2011	100
O2SI	020620-02	n-Hexane	1000	163378-27889	10-26-11B	11/14/2011	200
O2SI	020546-02	Heptane	1000	169174-29248	10-28-11G	11/14/2011	200
J&T Brand		Purge & Trap MeOH		K14E06-00556	10/27/2011	6/8/2012	3300
10-28-11AA							
50ug/ml Vol Work Std #2							
Exp:11/04/11							
Supplier	ID #	ID	ug/ml				
O2SI	121020-05	HSL'S-Ketone Solution	2000	169173-28307	10-12-11B	11/14/2011	100
J&T Brand		Purge & Trap MeOH		K14E06-00556	10/27/2011	10/14/2012	3900
		10-28-11AB	Exp:	11/4/2011			
		5ug/ml Vol Work Std #9					
		SOURCES	Lot	APPL Code	APPL Exp Date	ul	
		50ug/ml Vol Work Std #7		10-28-11X	10/31/2011	200	
		50ug/ml Vol Work Std #8		10-28-11Z	10/31/2011	200	
		J&T Brand		10/6/2011	6/8/2012	1600	
		10-28-11AC	Exp:	11/4/2011			
		5ug/ml Vol Work Std #10					
		SOURCES	Lot	APPL Code	APPL Exp Date	ul	
		50ug/ml Vol Work Std #1		10-28-11Y	10/31/2011	200	
		J&T Brand		10/27/2011	6/8/2012	1800	

GCMS STANDARD PREPARATION BOOK # _____ PAGE # 4

		10-28-11AD	Exp:	11/5/2011			
		5ug/ml Vol Work Std #12					
		SOURCE	Lot	APPL Code	APPL Exp Date	ul	
		50ug/ml Vol Work Std #2		10-28-11AA	10/31/2011	200	
		J&T Brand		10/27/2011	6/8/2012	1800	
10-28-11AE							
50ug/ml 8260 Surrogate			Conc.		Date	Exp.	
Exp:11/04/11			ug/ml	Lot #	Code	Date	uL
02SI	120002-01	8260B Surr Solution	2000	173249-28847	10-28-11H	11/14/2011	100
J&T Brand		Purge & Trap MeOH		X14E05-00556	10/27/2011	6/8/2012	3900
10-28-11AF			Exp:	11/5/2011			
5.0ug/ml 8260 Surrogate			Lot	APPL Code	APPL Exp Date	ul	
		50ug/ml 8260 Surrogate		10-28-11AE	10/31/2011	200	
J&T Brand		Purge & Trap MeOH		X14E05-00556	10/27/2011	1800	
10-28-11AG							
250ug/ml TBA/IBA/Acetonitrile/Cyclohexanone/Acrolein/2-P						APPL	
Exp:11/04/11			Conc.		Date	Exp.	
Supplier	ID #		ug/ml	Lot #	Code	Date	uL
02SI	120166-01	Volatile Mix 4-3	2000	178651-29510	10-28-11I	12/17/2011	500
02SI	020229-09	Acrolein	10000	179961-29661	10-19-11H	11/21/2011	100
J&T Brand		Purge & Trap MeOH		X14E05-00556	10/27/2011	6/8/2012	3400

10/28/11 RS.

10/30/11 RS.

10/30/11 RS.

10/30/11 RS.

10/31/11 RS.

NOTEBOOK INSERT LABEL

Gasoline
Lot: LB82077 EXP: FEB/2014 STORAGE: ROOM TEMP. 47616-U 1 x 1ml

DATE RECEIVED:_____

SUPELCO
595 North Harrison Road • Bellefonte, PA
16823-0048 USA • Phone 814-359-3441

RS.

A -

STANDARD TRANSFER LABEL

Date of Preparation:
Reference Number:
Description: Gasoline
Lot #: LB82077 - 29133
Rec: 8/4/11 MFR exp. 02/28/14
Exp. Date: EXP: FEB/2014
Storage: ROOM TEMP.

B -

REST
Cat# 30205
Unleaded gasoline composite
Lot #: A076842 - 29141
Rec: 8/4/11 MFR exp. 10/31/17
Unleaded Gasoline Composite Standard
50000 ug/mL each in P&T Methanol
Lot# A076842 Exp.Date: 10/2017 Store: Freezer
Restek Corporation - 110 Benner Circle - Bellefonte, PA 16823

RS.

10/30/11C							
2000ug/ml Gasoline						APPL	
			Conc.		Date	Exp.	
Supplier	ID #		ug/ml	Lot #	Code	Date	uL
Supelco	LB82077	Gasoline	20,000	LB82077-29133	10-30-11A	11/2/2012	200
J&T Brand		Purge & Trap MeOH		X14E06-00556	10/27/2011	3/2/2012	18000
10/30/11D						APPL	
2000ug/ml Unleaded Gasoline						Exp.	
			Conc.		Date	Date	uL
Supplier	ID #		ug/ml	Lot #	Code	11/30/2012	80
Supelco	30205	Unleaded Gasoline	50,000	A076842-29141	10-30-11B	3/2/2012	31920
J&T Brand		Purge & Trap MeOH		X14E06-00556	10/27/2011		

GC/MS STANDARD PREPARATION BOOK # _____ PAGE # _____

E

Custom VOC Mix, 16-4, 100 mg/L, 4 x 1 mL
232725-03-4PAK
Lot # 162917 | Storage ≤ -10 Degrees | Expiry 8/11/12
Solvent: P/T Methanol

Custom VOC Mix 16-4
Lot #: 162917 - 27029
Rec: 8/13/10 MFR exp. 08/11/12

CHICO RS 12/07/11.

Gasoline Curve Preparation for 100mL Purge (water)-CHICO			
	Expiration Date:	1µg/mL Gasoline	10/31/2011 Final Vol
Date	Conc.	10-30-11C	w/P&T H2O
Code	µg/L	Exp:12-27-12	mL
10-30-11F	20	1	100
10-30-11G	60	2.5	100
10-30-11H	100	5	100
10-30-11I	300	15	100
10-30-11J	600	30	100
10-30-11K	800	40	100
10-30-11L	1000	50	100

RS.

Volatile Standard Curve Preparation for 10mL Purge (8260 water)-CHICO												
	Expiration Date:		10/31/2011									
		5µg/mL Vol Std #9	5µg/mL Surr	50µg/mL Vol Std #7	50µg/mL Vol Std #8	50µg/mL Surr	5µg/mL Vol Std #10	50µg/mL Vol Std #1	50µg/mL Vol Std #2	5µg/mL Vol Std #12		
Date	Conc.	10-28-11O	10-28-11S	10-28-11K	10-28-11M	10-28-11R	10-28-11P	10-28-11L	10-28-11N	10-28-11Q		
Code	µg/L	Exp:11-04-11	Exp:11-04-11	Exp:11-04-11	Exp:11-04-11	Exp:11-04-11	Exp:11-04-11	Exp:11-04-11	Exp:11-04-11	Exp:11-04-11		
10-30-11I	0.3	3	6	n/a	n/a	n/a	3	n/a	n/a	3		
10-30-11J	0.5	5	10	n/a	n/a	n/a	5	n/a	n/a	5		
10-30-11K	1	10	20	n/a	n/a	n/a	10	n/a	n/a	10		
10-30-11L	2	20	40	n/a	n/a	n/a	20	n/a	n/a	20		
10-30-11M	5	n/a	n/a	5	5	10	n/a	5	5	n/a		
10-30-11N	10	n/a	n/a	10	10	25	n/a	10	20	n/a		
10-30-11O	20	n/a	n/a	20	20	40	n/a	20	20	n/a		
10-30-11P	40	n/a	n/a	40	40	80	n/a	40	40	n/a		
10-30-11Q	100	n/a	n/a	100	100	n/a	n/a	100	100	n/a		

RS

290µg/mL TAPD 10-28-11T	Final Vol w/P&T H2O
Exp:11-04-11	mL
3	50
5	50
10	50
15	50
20	50
25	50
30	50
35	50
40	50

Volatile Standard Curve Preparation for 10mL Purge (8260 water)-MAX												
	Expiration Date:		11/1/2011									
		5µg/mL Vol Std #9	5µg/mL Surr	50µg/mL Vol Std #7	50µg/mL Vol Std #8	50µg/mL Surr	5µg/mL Vol Std #10	50µg/mL Vol Std #1	50µg/mL Vol Std #2	5µg/mL Vol Std #12		
Date	Conc.	10-28-11O	10-28-11S	10-28-11K	10-28-11M	10-28-11R	10-28-11P	10-28-11L	10-28-11N	10-28-11Q		
Code	µg/L	Exp:11-04-11	Exp:11-04-11	Exp:11-04-11	Exp:11-04-11	Exp:11-04-11	Exp:11-04-11	Exp:11-04-11	Exp:11-04-11	Exp:11-04-11		
10-31-11A	0.3	3	6	n/a	n/a	n/a	3	n/a	n/a	3		
10-31-11B	0.5	5	10	n/a	n/a	n/a	5	n/a	n/a	5		
10-31-11C	1	10	20	n/a	n/a	n/a	10	n/a	n/a	10		
10-31-11D	2	20	40	n/a	n/a	n/a	20	n/a	n/a	20		
10-31-11E	5	n/a	n/a	5	5	10	n/a	5	5	n/a		
10-31-11F	10	n/a	n/a	10	10	25	n/a	10	20	n/a		
10-31-11G	20	n/a	n/a	20	20	40	n/a	20	20	n/a		
10-31-11H	40	n/a	n/a	40	40	80	n/a	40	40	n/a		
10-31-11I	100	n/a	n/a	100	100	n/a	n/a	100	100	n/a		

10/31/11 RS

RS

290µg/mL TAPD 10-28-11T	Final Vol w/P&T H2O
Exp:11-04-11	mL
3	50
5	50
10	50
15	50
20	50
25	50
30	50
35	50
40	50

Injection Log

Directory: M:\CHICO\DATA\C111030\

Line	Vial	FileName	Multiplier	SampleName	Misc Info	Injected
1	1	1030C13W.[	1	20ug/mL BFB STD 10-19-11	Water 2ul	30 Oct 11 22:01
2	1	1030C15W.[	1	Voc Std 10-30-11@0.3ug/L	Water 10mLw/ IS:10-30-11	30 Oct 11 23:28
3	1	1030C16W.[	1	Voc Std 10-30-11@0.5ug/L	Water 10mLw/ IS:10-30-11	31 Oct 11 00:11
4	1	1030C17W.[	1	Voc Std 10-30-11@1.0ug/L	Water 10mLw/ IS:10-30-11	31 Oct 11 00:54
5	1	1030C18W.[	1	Voc Std 10-30-11@2.0ug/L	Water 10mLw/ IS:10-30-11	31 Oct 11 1:37
6	1	1030C19W.[	1	Voc Std 10-30-11@5.0ug/L	Water 10mLw/ IS:10-30-11	31 Oct 11 2:20
7	1	1030C20W.[	1	Voc Std 10-30-11@10ug/L	Water 10mLw/ IS:10-30-11	31 Oct 11 3:03
8	1	1030C21W.[	1	Voc Std 10-30-11@20ug/L	Water 10mLw/ IS:10-30-11	31 Oct 11 3:46
9	1	1030C22W.[	1	Voc Std 10-30-11@40ug/L	Water 10mLw/ IS:10-30-11	31 Oct 11 4:29
10	1	1030C23W.[	1	Voc Std 10-30-11@100ug/L	Water 10mLw/ IS:10-30-11	31 Oct 11 5:12
11	1	1030C26W.[	1	20ug/mL BFB STD 10-19-11	Water 2ul	31 Oct 11 7:21
12	1	1030C27W.[	1	Voc Std 10-30-11@10ug/L	Water 10mLw/ IS&S:10-30/1	31 Oct 11 8:05
13	1	1030C28W.[	1	111030A LCS-1WC (SS)	Water 10mLw/ IS&S:10-30/1	31 Oct 11 8:48
14	1	1030C29W.[	1	GAS 300ug/L (SS)	Water 10mLw/ IS&S:10-30/1	31 Oct 11 9:31
15	1	1030C30W.[	1	GAS 300ug/L CCV-1WC	Water 10mLw/ IS&S:10-30/1	31 Oct 11 10:14
16	1	1030C31W.[	1	GAS 300ug/L LCS-1WC	Water 10mLw/ IS&S:10-30/1	31 Oct 11 10:57
17	1	1030C34W.[	1	111030A BLk-1WC	Water 10mLw/ IS&S:10-30/1	31 Oct 11 13:02
18	1	1030C35W.[	1	AY49335W01	Water 10mLw/ IS&S:10-30/1	31 Oct 11 13:39
19	1	1030C36W.[	1	AY49333W04	Water 10mLw/ IS&S:10-30/1	31 Oct 11 14:16
20	1	1030C37W.[	1	AY49334W13	Water 10mLw/ IS&S:10-30/1	31 Oct 11 14:53
21	1	1030C38W.[	1	AY49336W04	Water 10mLw/ IS&S:10-30/1	31 Oct 11 15:31
22	1	1030C39W.[	1	AY49334W141516 MS-1WC (VOC)	Water 10mLw/ IS&S:10-30/1	31 Oct 11 16:08
23	1	1030C40W.[	1	AY49334W141516 MSD-1WC (VOC)	Water 10mLw/ IS&S:10-30/1	31 Oct 11 16:45
24	1	1030C41W.[	1	AY49334W161718 MS-1WC (GAS)	Water 10mLw/ IS&S:10-30/1	31 Oct 11 17:22
25	1	1030C42W.[	1	AY49334W161718 MSD-1WC (GAS)	Water 10mLw/ IS&S:10-30/1	31 Oct 11 17:59

METALS

METALS
QC Summary

APPL, INC.

METALS BLANK

APPL Inc.
908 North Temperance Avenue
Clovis, CA 93611

Method	Analyte	Result	LOQ	LOD	DL	Units	Prep Date	Analysis Date	QC Group
6020	LEAD (PB) (DISSOL	0.19 J	0.5	0.22	0.11	ug/L	11/10/11	11/11/11	#602D-111110A-AY49334

J = Estimated value.

Laboratory Control Spike Recovery
METALS

APPL Inc.
908 North Temperance Avenue
Clovis, CA 93611

Method	Compound Name	Spike Level ug/L	SPK Result ug/L	SPK % Recovery	Recovery Limits	Extract Date	Analysis Date	QC Group
6020	LEAD (PB) (DISSOLVED)	50.0	50.0	100	80-120	11/10/2011	1/11/2011	#602D-111110A-AY49334

Comments: _____

Matrix Spike Recoveries

METALS

APPL ID: **111110W-49334 MS - 161255**

Sample ID: AY49334

Client ID: ES047

APPL Inc.
908 North Temperance Avenue
Clovis, CA 93611

Method	Compound Name	Spike Lvl ug/L	Matrix Res ug/L	SPK Res ug/L	DUP Res ug/L	SPK % Recovery	DUP % Recovery	RPD	RPD Max	Recovery Limits	Extract Date-Spk	Analysis Date-Spk	Extract Date-Dup	Analysis Date-Dup	QC Group	QC Sample
6020	LEAD (PB) (DISSOLVE	50.0	ND	49.5	48.8	99.0	97.6	1.4	20	80-120	1/10/2011	1/11/2011	1/10/2011	1/11/2011	161255	AY49334

Comments: _____

METALS
Sample Data

APPL, INC.

Metals Analysis

Environet, Inc.
650 Iwilei Rd, #204
Honolulu, HI 96817

APPL Inc.
908 North Temperance Avenue
Clovis, CA 93611

Attn: Stacy Fineran
Project: RED HILL/1022-024
Sample ID: ES046
Sample Collection Date: 10/24/2011

ARF: 66102
APPL ID: AY49333

Method	Analyte	Result	LOQ	LOD	DL	Units	DF	Prep Date	Analysis Date
6020	LEAD (PB) (DISSOLVED)	0.22 U	0.5	0.22	0.11	ug/L	1	11/10/2011	11/11/2011

Printed: 11/15/2011 12:57:19 PM

APPL-F1-SC-NoMC-REG MDLs

Sample QC Report

Data File: C:\ICPCHEM\1\DATA\11K11100.B\041SMPL.D\041SMPL.D#
Date Acquired: Nov 11 2011 04:00 pm
Operator: NBS
Sample Name: AY49333W13
Misc Info: 111110A-3015
Vial Number: 3202
Current Method: C:\ICPCHEM\1\METHODS\62A1111A.M
Calibration File: C:\ICPCHEM\1\CALIB\62A1111A.C
Last Cal Update: Nov 11 2011 12:36 pm
Sample Type: Sample
Prep Dil Factor: 1.11
Total Dil Factor: 1.11

QC Elements

Element	Conc.	Corr. Conc.	RSD(%)	High Limit	Flag
7 (Li)	-------- ug/l	#VALUE!	--------	0	
9 Be	-0.01 ug/l	-0.01	5.25	1000	
11 B	74.87 ug/l	83.18	0.56	1000	
23 Na	63250.00 ug/l	70270.75	0.40	25000	>Cal
24 Mg	23790.00 ug/l	26430.69	0.63	50000	
27 Al	35.06 ug/l	38.95	3.68	20000	
39 K	2658.00 ug/l	2953.04	0.91	20000	
44 Ca	12970.00 ug/l	14409.67	0.32	50000	
47 Ti	1.59 ug/l	1.77	2.95	1000	
51 V	1.07 ug/l	1.18	4.80	1000	
52 Cr	0.15 ug/l	0.16	6.59	1000	
55 Mn	1421.00 ug/l	1578.73	0.18	1000	>Cal
56 Fe	849.00 ug/l	943.24	0.37	20000	
59 Co	0.20 ug/l	0.22	4.00	1000	
60 Ni	1.09 ug/l	1.21	7.81	1000	
63 Cu	-0.35 ug/l	-0.39	4.23	1000	
65 Cu	-0.35 ug/l	-0.38	7.78	1000	
66 Zn	1.22 ug/l	1.36	1.97	1000	
75 As	0.07 ug/l	0.08	32.39	1000	
78 Se	0.04 ug/l	0.04	27.66	1000	
78 Se	0.16 ug/l	0.18	40.54	1000	
88 Sr	115.00 ug/l	127.77	1.27	1000	
88 Sr	117.90 ug/l	130.99	0.51	1000	
95 Mo	0.48 ug/l	0.54	2.21	1000	
106 (Cd)	-------- ug/l	#VALUE!	--------	#######	
107 Ag	0.00 ug/l	0.00	108.51	500	
108 (Cd)	-------- ug/l	#VALUE!	--------	#######	
111 Cd	0.01 ug/l	0.02	47.56	1000	
118 Sn	0.11 ug/l	0.13	17.45	1000	
121 Sb	0.07 ug/l	0.08	13.55	1000	
137 Ba	23.17 ug/l	25.74	1.06	1000	
205 Tl	0.01 ug/l	0.01	19.54	1000	
206 (Pb)	-------- ug/l	#VALUE!	--------	#######	
207 (Pb)	-------- ug/l	#VALUE!	--------	#######	
208 Pb	-0.16 ug/l	-0.18	5.22	1000	

ISTD Elements

Element	CPS Mean	RSD(%)	Ref Value	Rec(%)	QC Range(%)	Flag
6 Li	3155449.80	0.29	2775704.50	113.7	70 - 120	
45 Sc	684547.81	0.76	500780.41	136.7	70 - 120	IS Fai
45 Sc	107493.14	1.57	95494.08	112.6	70 - 120	
45 Sc	2252708.50	1.98	1460980.80	154.2	70 - 120	IS Fai
72 Ge	105205.95	0.22	96219.04	109.3	70 - 120	
72 Ge	48357.57	0.66	43611.78	110.9	70 - 120	
72 Ge	228774.88	0.47	213204.63	107.3	70 - 120	
115 In	1463890.90	0.38	1381264.00	106.0	70 - 120	
159 Tb	2015104.00	0.80	1843940.90	109.3	70 - 120	
165 Ho	2022052.00	0.63	1844184.90	109.6	70 - 120	

ISTD Ref File : C:\ICPCHEM\1\DATA\11K11100.B\004CALB.D\004CALB.D#

2 :Element Failures 0 :Max. Number of Failures Allowed
2 :ISTD Failures 0 :Max. Number of ISTD Failures Allowed

Data Results:
Analytes: Fail
ISTD: Fail

Metals Analysis

Environet, Inc.
650 Iwilei Rd, #204
Honolulu, HI 96817

Attn: Stacy Fineran

Project: RED HILL/1022-024

Sample ID: ES047

Sample Collection Date: 10/24/2011

APPL Inc.
908 North Temperance Avenue
Clovis, CA 93611

ARF: 66102

APPL ID: AY49334

Method	Analyte	Result	LOQ	LOD	DL	Units	DF	Prep Date	Analysis Date
6020	LEAD (PB) (DISSOLVED)	0.22U	0.5	0.22	0.11	ug/L	1	11/10/2011	11/11/2011

Printed: 11/15/2011 12:57:19 PM

APPL-F1-SC-NoMC-REG MDLs

Sample QC Report

Data File: C:\ICPCHEM\1\DATA\11K1100.B\042SMPL.D\042SMPL.D#
Date Acquired: Nov 11 2011 04:06 pm
Operator: NBS
Sample Name: AY49334W51
Misc Info: 111110A-3015
Vial Number: 3203
Current Method: C:\ICPCHEM\1\METHODS\62A1111A.M
Calibration File: C:\ICPCHEM\1\CALIB\62A1111A.C
Last Cal Update: Nov 11 2011 12:36 pm
Sample Type: Sample
Prep Dil Factor: 1.11
Total Dil Factor: 1.11

QC Elements

Element	Conc.		Corr. Conc.	RSD(%)	High Limit	Flag
7 (Li)	--------	ug/l	#VALUE!	--------	0	
9 Be	-0.01	ug/l	-0.01	2.99	1000	
11 B	77.21	ug/l	85.78	0.37	1000	
23 Na	62580.00	ug/l	69526.38	1.20	25000	>Cal
24 Mg	23840.00	ug/l	26486.24	0.85	50000	
27 Al	10.66	ug/l	11.84	1.02	20000	
39 K	2644.00	ug/l	2937.48	0.98	20000	
44 Ca	12880.00	ug/l	14309.68	1.08	50000	
47 Ti	1.96	ug/l	2.17	54.12	1000	
51 V	1.12	ug/l	1.24	3.35	1000	
52 Cr	0.18	ug/l	0.20	5.89	1000	
55 Mn	1422.00	ug/l	1579.84	0.97	1000	>Cal
56 Fe	862.60	ug/l	958.35	1.25	20000	
59 Co	0.26	ug/l	0.29	3.17	1000	
60 Ni	1.23	ug/l	1.37	2.27	1000	
63 Cu	-0.28	ug/l	-0.31	4.35	1000	
65 Cu	-0.29	ug/l	-0.32	5.75	1000	
66 Zn	5.31	ug/l	5.90	2.33	1000	
75 As	0.09	ug/l	0.10	35.61	1000	
78 Se	0.04	ug/l	0.04	18.02	1000	
78 Se	0.13	ug/l	0.15	43.74	1000	
88 Sr	114.00	ug/l	126.65	0.20	1000	
88 Sr	116.30	ug/l	129.21	0.82	1000	
95 Mo	0.49	ug/l	0.54	2.47	1000	
106 (Cd)	--------	ug/l	#VALUE!	--------	#######	
107 Ag	0.00	ug/l	0.00	9066.00	500	
108 (Cd)	--------	ug/l	#VALUE!	--------	#######	
111 Cd	0.00	ug/l	0.00	326.50	1000	
118 Sn	0.09	ug/l	0.11	17.30	1000	
121 Sb	0.05	ug/l	0.05	7.46	1000	
137 Ba	24.70	ug/l	27.44	0.69	1000	
205 Tl	0.01	ug/l	0.01	3.49	1000	
206 (Pb)	--------	ug/l	#VALUE!	--------	#######	
207 (Pb)	--------	ug/l	#VALUE!	--------	#######	
208 Pb	-0.05	ug/l	-0.06	14.15	1000	

ISTD Elements

Element	CPS Mean	RSD(%)	Ref Value	Rec(%)	QC Range(%)	Flag
6 Li	3186521.00	0.78	2775704.50	114.8	70 - 120	
45 Sc	672635.13	1.27	500780.41	134.3	70 - 120	IS Fai
45 Sc	107878.18	1.34	95494.08	113.0	70 - 120	
45 Sc	2212596.80	1.01	1460980.80	151.4	70 - 120	IS Fai
72 Ge	106314.52	1.16	96219.04	110.5	70 - 120	
72 Ge	48500.91	1.75	43611.78	111.2	70 - 120	
72 Ge	231305.58	0.13	213204.63	108.5	70 - 120	
115 In	1470644.80	1.61	1381264.00	106.5	70 - 120	
159 Tb	2036736.10	0.19	1843940.90	110.5	70 - 120	
165 Ho	2036137.40	0.55	1844184.90	110.4	70 - 120	

ISTD Ref File : C:\ICPCHEM\1\DATA\11K1100.B\004CALB.D\004CALB.D#

2 :Element Failures 0 :Max. Number of Failures Allowed
2 :ISTD Failures 0 :Max. Number of ISTD Failures Allowed

Data Results:
Analytes: Fail
ISTD: Fail

Metals Analysis

Environet, Inc.
650 Iwilei Rd, #204
Honolulu, HI 96817

APPL Inc.
908 North Temperance Avenue
Clovis, CA 93611

Attn: Stacy Fineran
Project: RED HILL/1022-024
Sample ID: ES049
Sample Collection Date: 10/24/2011

ARF: 66102
APPL ID: AY49336

Method	Analyte	Result	LOQ	LOD	DL	Units	DF	Prep Date	Analysis Date
6020	LEAD (PB) (DISSOLVED)	0.22J	0.5	0.22	0.11	ug/L	1	11/10/2011	11/11/2011

J = Estimated value.

Printed: 11/15/2011 12:57:19 PM

APPL-F1-SC-NoMC-REG MDLs

Sample QC Report

Data File: C:\ICPCHEM\1\DATA\11K11100.B\052SMPL.D\052SMPL.D#
Date Acquired: Nov 11 2011 05:06 pm
Operator: NBS
Sample Name: AY49336W13
Misc Info: 111110A-3015
Vial Number: 3208
Current Method: C:\ICPCHEM\1\METHODS\62A1111A.M
Calibration File: C:\ICPCHEM\1\CALIB\62A1111A.C
Last Cal Update: Nov 11 2011 12:36 pm
Sample Type: Sample
Prep Dil Factor: 1.11
Total Dil Factor: 1.11

QC Elements

Element	Conc.	Corr. Conc.	RSD(%)	High Limit	Flag
7 (Li)	-------- ug/l	#VALUE!	--------	0	
9 Be	-0.01 ug/l	-0.01	2.49	1000	
11 B	119.20 ug/l	132.43	0.84	1000	
23 Na	79710.00 ug/l	88557.81	0.80	25000	>Cal
24 Mg	26180.00 ug/l	29085.98	1.28	50000	
27 Al	28.51 ug/l	31.67	0.66	20000	
39 K	3332.00 ug/l	3701.85	0.44	20000	
44 Ca	22270.00 ug/l	24741.97	1.18	50000	
47 Ti	0.71 ug/l	0.79	6.25	1000	
51 V	15.17 ug/l	16.85	0.80	1000	
52 Cr	0.80 ug/l	0.89	2.21	1000	
55 Mn	106.10 ug/l	117.88	0.65	1000	
56 Fe	83.80 ug/l	93.10	0.12	20000	
59 Co	1.02 ug/l	1.13	0.83	1000	
60 Ni	3.41 ug/l	3.79	1.43	1000	
63 Cu	0.33 ug/l	0.36	4.20	1000	
65 Cu	0.33 ug/l	0.37	5.05	1000	
66 Zn	4.07 ug/l	4.52	2.37	1000	
75 As	0.30 ug/l	0.34	16.47	1000	
78 Se	0.13 ug/l	0.14	5.92	1000	
78 Se	0.15 ug/l	0.16	65.10	1000	
88 Sr	150.20 ug/l	166.87	1.25	1000	
88 Sr	151.90 ug/l	168.76	0.26	1000	
95 Mo	19.64 ug/l	21.82	0.46	1000	
106 (Cd)	-------- ug/l	#VALUE!	--------	#######	
107 Ag	0.06 ug/l	0.06	16.30	500	
108 (Cd)	-------- ug/l	#VALUE!	--------	#######	
111 Cd	0.01 ug/l	0.01	93.49	1000	
118 Sn	0.18 ug/l	0.20	15.54	1000	
121 Sb	0.74 ug/l	0.83	4.50	1000	
137 Ba	12.11 ug/l	13.45	0.77	1000	
205 Tl	0.02 ug/l	0.02	13.88	1000	
206 (Pb)	-------- ug/l	#VALUE!	--------	#######	
207 (Pb)	-------- ug/l	#VALUE!	--------	#######	
208 Pb	0.20 ug/l	0.22	6.86	1000	

ISTD Elements

Element	CPS Mean	RSD(%)	Ref Value	Rec(%)	QC Range(%)	Flag
6 Li	3041975.00	0.97	2775704.50	109.6	70 - 120	
45 Sc	642764.69	0.96	500780.41	128.4	70 - 120	IS Fai
45 Sc	102731.88	1.25	95494.08	107.6	70 - 120	
45 Sc	2133139.30	1.51	1460980.80	146.0	70 - 120	IS Fai
72 Ge	102652.80	0.93	96219.04	106.7	70 - 120	
72 Ge	46837.37	1.69	43611.78	107.4	70 - 120	
72 Ge	230432.05	0.56	213204.63	108.1	70 - 120	
115 In	1482746.90	1.89	1381264.00	107.3	70 - 120	
159 Tb	2029492.10	1.13	1843940.90	110.1	70 - 120	
165 Ho	2027253.00	1.23	1844184.90	109.9	70 - 120	

ISTD Ref File : C:\ICPCHEM\1\DATA\11K11100.B\004CALB.D\004CALB.D#

1 :Element Failures 0 :Max. Number of Failures Allowed
2 :ISTD Failures 0 :Max. Number of ISTD Failures Allowed

Data Results:
Analytes: Fail
ISTD: Fail

METALS
Calibration Data

APPL, INC.

A.P.P.L. INC.

2A

INITIAL AND CONTINUING CALIBRATION VERIFICATION

Lab Name: A.P.P.L. INC. Contract: Environet, Inc.

ARF No: 66102 SDG: 66102

Initial Calibration Source: CPI

Continuing Calibration Source: Environmental Express

Analysis Date: 11/11/2011 Concentration Units: ug/L

Analyte	Initial Calibration			Continuing Calibration						M
	True	Found 12:39	%R(1)	True CCV1	Found 13:03	%R(1)	True CCV1	Found 13:33	%R(1)	
Lead (Pb)	100	106.3	106	50	50.31	101	50	50.34	101	P

(1) Control Limits: Metals 90-110

ILM02.0

A.P.P.L. INC.

2A

INITIAL AND CONTINUING CALIBRATION VERIFICATION

Lab Name: A.P.P.L. INC. Contract: Environet, Inc.

ARF No: 66102 SDG: 66102

Initial Calibration Source: CPI

Continuing Calibration Source: Environmental Express

Analysis Date: 11/11/2011 Concentration Units: ug/L

Analyte	Initial Calibration			Continuing Calibration						M
	True	Found 12:39	%R(1)	True CCV1	Found 15:05	%R(1)	True CCV1	Found 16:30	%R(1)	
Lead (Pb)	100	106.3	106	50	49.96	99.9	50	50.41	101	P

(1) Control Limits: Metals 90-110

ILM02.0

A.P.P.L. INC.

2A

INITIAL AND CONTINUING CALIBRATION VERIFICATION

Lab Name: A.P.P.L. INC. Contract: Environet, Inc.

ARF No: 66102 SDG: 66102

Initial Calibration Source: CPI

Continuing Calibration Source: Environmental Express

Analysis Date: 11/11/2011 Concentration Units: ug/L

Analyte	Initial Calibration			Continuing Calibration						M
	True	Found 12:39	%R(1)	True CCV1	Found 18:04	%R(1)	True	Found	%R(1)	
Lead (Pb)	100	106.3	106	50	48.04	96.1				P

(1) Control Limits: Metals 90-110

ILM02.0

A.P.P.L. INC.
3
BLANKS

Lab Name: A.P.P.L. INC. Contract: Environet, Inc.

ARF No.: 66102 SDG: 66102

Preparation Blank Matrix (soil/water): water

Preparation Blank Concentration Units (ug/L or mg/kg): ug/L

Analysis Date: 11/11/2011

Analyte	Initial Calibration Blank (ug/L)		Continuing Calibration Blank (ug/L)								Preparation Blank		M
		C	1	C	2	C	3	C		C			
	12:57		13:09		13:46		15:17		14:16				
Lead (Pb)	.50	U	.50	U	.50	U	.50	U	.19	J	P		

A.P.P.L. INC.
3
BLANKS

Lab Name: A.P.P.L. INC. Contract: Environet, Inc.

ARF No.: 66102 SDG: 66102

Preparation Blank Matrix (soil/water): water

Preparation Blank Concentration Units (ug/L or mg/kg): ug/L

Analysis Date: 11/11/2011

Analyte	Initial Calibration Blank (ug/L)		Continuing Calibration Blank (ug/L)						Preparation Blank		M
		C	1	C	2	C	3	C		C	
	12:57		16:42		18:16				14:16		
Lead (Pb)	.50	U	.50	U	.50	U			.19	J	P

A.P.P.L. INC.
4
ICP INTERFERENCE CHECK SAMPLE

Lab Name: A.P.P.L. INC. Contract: Environet, Inc.
ARF No.: 66102 SDG: 66102
ICP ID Number: Optimus ICS Source: Environmental Express

Analysis Date: 11/11/2011 Concentration Units: ug/L

Analyte	True		Initial Found		
	Sol A	Sol AB	Sol A 13:15	Sol AB 13:21	%R(1)
Lead (Pb)		500	3.499	502	100

(1) Control Limits: Metals 80-120

A.P.P.L. INC.

5B

POST DIGEST SPIKE SAMPLE RECOVERY

CLIENT SAMPLE NO.

ES047

Lab Name: A.P.P.L. INC. Contract: Environet, Inc.

ARF No.: 66102 SDG: 66102

Analysis Date: 11/11/2011 Concentration Units: ug/L

Analyte	Control Limit %R	Spiked Sample Result (SSR)	C	Sample Result (SR)	C	Spike Added (SA)	%R	Q	M
Lead (Pb)	75-125	259.629		-0.0610389		277.500	93.6		

Comments:

11/11/201116:06 AY49334W51

11/11/201116:54 AY49334W51-A

Sample QC Report

Data File: C:\ICPCHEM\1\DATA\11K1110O.B\050SMPL.D\050SMPL.D#
Date Acquired: Nov 11 2011 04:54 pm
Operator: NBS
Sample Name: AY49334W51-A
Misc Info: 111110A-3015
Vial Number: 3206
Current Method: C:\ICPCHEM\1\METHODS\62A1111A.M
Calibration File: C:\ICPCHEM\1\CALIB\62A1111A.C
Last Cal Update: Nov 11 2011 12:36 pm
Sample Type: Sample
Prep Dil Factor: 1.11
Total Dil Factor: 1.11

QC Elements

Element	Conc.		Corr. Conc.	RSD(%)	High Limit	Flag
7 (Li)	--------	ug/l	#VALUE!	--------	0	
9 Be	27.12	ug/l	30.13	1.56	1000	
11 B	219.90	ug/l	244.31	1.71	1000	
23 Na	81860.00	ug/l	90946.46	1.33	25000	>Cal
24 Mg	43910.00	ug/l	48784.01	1.40	50000	
27 Al	1865.00	ug/l	2072.02	1.45	20000	
39 K	7013.00	ug/l	7791.44	0.83	20000	
44 Ca	35340.00	ug/l	39262.74	0.74	50000	
47 Ti	223.10	ug/l	247.86	0.84	1000	
51 V	240.60	ug/l	267.31	0.69	1000	
52 Cr	236.20	ug/l	262.42	1.19	1000	
55 Mn	1601.00	ug/l	1778.71	1.31	1000	>Cal
56 Fe	1722.00	ug/l	1913.14	1.16	20000	
59 Co	239.30	ug/l	265.86	1.18	1000	
60 Ni	220.50	ug/l	244.98	0.57	1000	
63 Cu	217.50	ug/l	241.64	0.86	1000	
65 Cu	218.50	ug/l	242.75	0.57	1000	
66 Zn	414.80	ug/l	460.84	0.80	1000	
75 As	221.10	ug/l	245.64	0.31	1000	
78 Se	198.20	ug/l	220.20	2.61	1000	
78 Se	205.70	ug/l	228.53	1.47	1000	
88 Sr	383.00	ug/l	425.51	0.35	1000	
88 Sr	358.90	ug/l	398.74	1.23	1000	
95 Mo	242.00	ug/l	268.86	1.21	1000	
106 (Cd)	--------	ug/l	#VALUE!	--------	#######	
107 Ag	68.49	ug/l	76.09	4.03	500	
108 (Cd)	--------	ug/l	#VALUE!	--------	#######	
111 Cd	44.30	ug/l	49.22	1.31	1000	
118 Sn	271.00	ug/l	301.08	1.25	1000	
121 Sb	243.60	ug/l	270.64	0.29	1000	
137 Ba	267.20	ug/l	296.86	0.54	1000	
205 Tl	227.70	ug/l	252.97	0.11	1000	
206 (Pb)	--------	ug/l	#VALUE!	--------	#######	
207 (Pb)	--------	ug/l	#VALUE!	--------	#######	
208 Pb	233.90	ug/l	259.86	1.71	1000	

ISTD Elements

Element	CPS Mean	RSD(%)	Ref Value	Rec(%)	QC Range(%)	Flag
6 Li	2876337.50	0.73	2775704.50	103.6	70 - 120	
45 Sc	648562.19	0.34	500780.41	129.5	70 - 120	IS Fai
45 Sc	102181.34	0.75	95494.08	107.0	70 - 120	
45 Sc	2157529.50	1.05	1460980.80	147.7	70 - 120	IS Fai
72 Ge	102359.33	0.69	96219.04	106.4	70 - 120	
72 Ge	46173.78	0.92	43611.78	105.9	70 - 120	
72 Ge	223471.42	0.75	213204.63	104.8	70 - 120	
115 In	1440331.40	0.34	1381264.00	104.3	70 - 120	
159 Tb	1995111.90	0.82	1843940.90	108.2	70 - 120	
165 Ho	1988682.90	0.53	1844184.90	107.8	70 - 120	

ISTD Ref File : C:\ICPCHEM\1\DATA\11K1110O.B\004CALB.D\004CALB.D#

2 :Element Failures 0 :Max. Number of Failures Allowed
2 :ISTD Failures 0 :Max. Number of ISTD Failures Allowed

Data Results:
Analytes: Fail
ISTD: Fail

A.P.P.L. INC.

9

ICP SERIAL DILUTION

CLIENT SAMPLE NO.

ES047

Lab Name: A.P.P.L. INC. Contract: Environet, Inc.

ARF No.: 66102 SDG: 66102

Matrix: water

Analysis Date: 11/11/2011 Concentration Units: ug/L

Analyte	Initial Sample Result (I)	C	Serial Dilution Result (S)	C	%D	Q	M
Lead (Pb)	-0.0610389		-1.23988		NA		

Comments:

11/11/2011 16:06 AY49334W51

11/11/2011 17:00 AY49334W51-1/5

Sample QC Report

Data File: C:\ICPCHEM\1\DATA\11K11100.B\051SMPL.D\051SMPL.D#
Date Acquired: Nov 11 2011 05:00 pm
Operator: NBS
Sample Name: AY49334W51-1/5
Misc Info: 111110A-3015
Vial Number: 3207
Current Method: C:\ICPCHEM\1\METHODS\62A1111A.M
Calibration File: C:\ICPCHEM\1\CALIB\62A1111A.C
Last Cal Update: Nov 11 2011 12:36 pm
Sample Type: Sample
Prep Dil Factor: 5.56
Total Dil Factor: 5.56

QC Elements

Element	Conc.	Corr. Conc.	RSD(%)	High Limit	Flag
7 (Li)	-------- ug/l	#VALUE!	--------	0	
9 Be	-0.01 ug/l	-0.07	3.30	1000	
11 B	17.93 ug/l	99.62	0.55	1000	
23 Na	13240.00 ug/l	73561.44	1.12	25000	
24 Mg	5084.00 ug/l	28246.70	1.02	50000	
27 Al	1.95 ug/l	10.81	25.75	20000	
39 K	473.00 ug/l	2627.99	1.52	20000	
44 Ca	2644.00 ug/l	14690.06	1.59	50000	
47 Ti	0.34 ug/l	1.89	8.30	1000	
51 V	3.79 ug/l	21.04	2.49	1000	
52 Cr	0.14 ug/l	0.80	4.27	1000	
55 Mn	287.90 ug/l	1599.57	0.52	1000	
56 Fe	175.80 ug/l	976.74	0.50	20000	
59 Co	-0.18 ug/l	-1.02	3.42	1000	
60 Ni	0.22 ug/l	1.23	6.12	1000	
63 Cu	-0.54 ug/l	-3.02	2.33	1000	
65 Cu	-0.54 ug/l	-3.00	2.74	1000	
66 Zn	1.00 ug/l	5.58	4.28	1000	
75 As	0.92 ug/l	5.11	1.72	1000	
78 Se	0.16 ug/l	0.87	1.85	1000	
78 Se	0.30 ug/l	1.65	38.09	1000	
88 Sr	22.51 ug/l	125.07	1.55	1000	
88 Sr	20.24 ug/l	112.45	0.28	1000	
95 Mo	0.18 ug/l	1.01	8.72	1000	
106 (Cd)	-------- ug/l	#VALUE!	--------	########	
107 Ag	0.83 ug/l	4.59	0.55	500	
108 (Cd)	-------- ug/l	#VALUE!	--------	########	
111 Cd	0.01 ug/l	0.07	56.81	1000	
118 Sn	0.30 ug/l	1.65	5.07	1000	
121 Sb	1.56 ug/l	8.68	3.00	1000	
137 Ba	4.97 ug/l	27.59	0.72	1000	
205 Tl	0.01 ug/l	0.06	8.19	1000	
206 (Pb)	-------- ug/l	#VALUE!	--------	########	
207 (Pb)	-------- ug/l	#VALUE!	--------	########	
208 Pb	-0.22 ug/l	-1.24	2.31	1000	

ISTD Elements

Element	CPS Mean	RSD(%)	Ref Value	Rec(%)	QC Range(%)	Flag
6 Li	3013682.80	1.03	2775704.50	108.6	70 - 120	
45 Sc	545314.75	0.37	500780.41	108.9	70 - 120	
45 Sc	95160.34	0.89	95494.08	99.7	70 - 120	
45 Sc	1580825.90	0.15	1460980.80	108.2	70 - 120	
72 Ge	104611.55	0.65	96219.04	108.7	70 - 120	
72 Ge	45792.49	0.44	43611.78	105.0	70 - 120	
72 Ge	227516.67	0.25	213204.63	106.7	70 - 120	
115 In	1384767.60	0.50	1381264.00	100.3	70 - 120	
159 Tb	1892859.60	0.65	1843940.90	102.7	70 - 120	
165 Ho	1897757.40	0.79	1844184.90	102.9	70 - 120	

ISTD Ref File : C:\ICPCHEM\1\DATA\11K11100.B\004CALB.D\004CALB.D#

0 :Element Failures 0 :Max. Number of Failures Allowed
0 :ISTD Failures 0 :Max. Number of ISTD Failures Allowed

Data Results:
Analytes: Pass
ISTD: Pass

Calibration Blank QC Report

Data File:	C:\ICPCHEM\1\DATA\11K11100.B\004CAL
Date Acquired:	Nov 11 2011 12:08 pm
Operator:	NBS
Sample Name:	Calibration Blank
Misc Info:	
Vial Number:	1102
Current Method:	C:\ICPCHEM\1\METHODS\62A1111A.M
Calibration File:	C:\ICPCHEM\1\CALIB\62A1111A.C
Last Cal Update:	Nov 11 2011 12:06 pm
Sample Type:	CalBlk
Total Dil Factor:	1.00

QC&ISTD Elements

Element		CPS Mean		SD	RSD(%)
6	Li	2775705.00	A	31080.00	1.12
7	(Li)	152897.91	P	508.10	0.33
9	Be	164.45	P	15.75	9.58
11	B	9503.37	P	213.80	2.25
23	Na	81958.40	P	248.30	0.30
24	Mg	134.45	P	6.94	5.16
27	Al	111.12	P	16.78	15.10
39	K	60334.78	P	2276.00	3.77
44	Ca	384.84	P	48.59	12.63
45	Sc	500780.41	P	2032.00	0.41
45	Sc	95494.08	P	252.60	0.26
45	Sc	1460981.00	A	25510.00	1.75
47	Ti	4.89	P	0.77	15.75
51	V	3955.25	P	110.20	2.79
52	Cr	547.13	P	20.02	3.66
55	Mn	165.78	P	8.57	5.17
56	Fe	5746.57	P	137.00	2.38
59	Co	1492.99	P	62.44	4.18
60	Ni	69.78	P	22.72	32.56
63	Cu	2222.87	P	55.11	2.48
65	Cu	1076.95	P	27.98	2.60
66	Zn	207.12	P	12.10	5.84
72	Ge	96219.04	P	484.10	0.50
72	Ge	43611.78	P	490.40	1.12
72	Ge	213204.59	P	1657.00	0.78
75	As	266.34	P	7.21	2.71
78	Se	4.67	P	1.53	32.74
78	Se	30.00	P	1.16	3.85
88	Sr	48.89	P	8.39	17.16
88	Sr	188.90	P	11.71	6.20
95	Mo	111.12	P	22.69	20.42
106	(Cd)	31.11	P	10.18	32.72
107	Ag	35.56	P	13.47	37.88
108	(Cd)	27.78	P	5.09	18.33
111	Cd	0.12	P	4.33	3513.10
115	In	1381264.00	A	15790.00	1.14
118	Sn	495.58	P	60.50	12.21
121	Sb	323.35	P	35.28	10.91
137	Ba	91.12	P	13.47	14.78
159	Tb	1843941.00	A	33820.00	1.83
165	Ho	1844185.00	A	22050.00	1.20
205	Tl	78.89	P	5.09	6.45
206	(Pb)	1670.17	P	51.97	3.11
207	(Pb)	1455.69	P	79.06	5.43
208	Pb	6738.71	P	70.43	1.05

Calibration Standard QC Report

Data File: C:\ICPCHEM\1\DATA\11K11100.B\005CALS.D\005CALS.D#
Date Acquired: Nov 11 2011 12:14 pm
Operator: NBS
Sample Name: 111111 Standard 1
Misc Info:
Vial Number: 1103
Current Method: C:\ICPCHEM\1\METHODS\62A1111A.M
Calibration File: C:\ICPCHEM\1\CALIB\62A1111A.C
Last Cal Update: Nov 11 2011 12:12 pm
Sample Type: CalStd
Total Dil Factor: 1.00

QC&ISTD Elements

Element	CPS Mean		SD	RSD(%)	Cal Coef
6 Li	2934478.00	A	12100.00	0.41	0.0000
7 (Li)	160812.41	P	1266.00	0.79	0.0000
9 Be	1031.18	P	27.15	2.63	0.0000
11 B	10014.79	P	224.60	2.24	0.0000
23 Na	101764.30	P	5296.00	5.20	0.0000
24 Mg	2435.84	P	56.81	2.33	0.0000
27 Al	465.58	P	50.04	10.75	0.0000
39 K	63456.74	P	1758.00	2.77	0.0000
44 Ca	441.03	P	5.03	1.14	0.0000
45 Sc	483714.81	P	17820.00	3.68	0.0000
45 Sc	96706.18	P	602.60	0.62	0.0000
45 Sc	1494561.00	A	14240.00	0.95	0.0000
47 Ti	16.89	P	3.36	19.87	0.0000
51 V	4556.33	P	51.66	1.13	0.0000
52 Cr	876.48	P	32.73	3.73	0.0000
55 Mn	7451.77	P	52.30	0.70	0.0000
56 Fe	12699.44	P	213.90	1.68	0.0000
59 Co	1820.58	P	82.65	4.54	0.0000
60 Ni	166.23	P	12.10	7.28	0.0000
63 Cu	3334.65	P	61.70	1.85	0.0000
65 Cu	1647.67	P	94.43	5.73	0.0000
66 Zn	231.56	P	11.34	4.90	0.0000
72 Ge	93081.49	P	2181.00	2.34	0.0000
72 Ge	43620.24	P	387.20	0.89	0.0000
72 Ge	210910.70	P	1414.00	0.67	0.0000
75 As	300.78	P	7.07	2.35	0.0000
78 Se	21.00	P	2.60	12.40	0.0000
78 Se	30.33	P	6.33	20.88	0.0000
88 Sr	303.35	P	25.17	8.30	0.0000
88 Sr	1913.54	P	79.67	4.16	0.0000
95 Mo	385.58	P	18.36	4.76	0.0000
106 (Cd)	51.11	P	6.94	13.58	0.0000
107 Ag	447.80	P	37.47	8.37	0.0000
108 (Cd)	28.89	P	17.10	59.19	0.0000
111 Cd	182.07	P	18.49	10.16	0.0000
115 In	1383497.00	A	12980.00	0.94	0.0000
118 Sn	901.17	P	20.10	2.23	0.0000
121 Sb	988.96	P	26.95	2.73	0.0000
137 Ba	304.46	P	49.48	16.25	0.0000
159 Tb	1838841.00	A	19950.00	1.08	0.0000
165 Ho	1842078.00	A	20850.00	1.13	0.0000
205 Tl	1497.92	P	40.19	2.68	0.0000
206 (Pb)	2154.70	P	105.60	4.90	0.0000
207 (Pb)	1842.42	P	104.10	5.65	0.0000
208 Pb	8565.85	P	320.10	3.74	0.0000

ISTD Elements

Element	CPS Mean	RSD(%)	Ref Value	Rec(%)	QC Range(%)		Flag
6 Li	2934478.30	0.41	2775704.50	105.7	70 -	120	
45 Sc	483714.78	3.68	500780.41	96.6	70 -	120	
45 Sc	96706.18	0.62	95494.08	101.3	70 -	120	
45 Sc	1494561.00	0.95	1460980.80	102.3	70 -	120	
72 Ge	93081.49	2.34	96219.04	96.7	70 -	120	
72 Ge	43620.24	0.89	43611.78	100.0	70 -	120	
72 Ge	210910.72	0.67	213204.63	98.9	70 -	120	
115 In	1383496.90	0.94	1381264.00	100.2	70 -	120	
159 Tb	1838841.50	1.08	1843940.90	99.7	70 -	120	
165 Ho	1842078.10	1.13	1844184.90	99.9	70 -	120	

ISTD Ref File : C:\ICPCHEM\1\DATA\11K11100.B\004CALB.D\004CALB.D#

--- :Element Failures --- :Max. Number of Failures Allowsd
0 :ISTD Failures 0 :Max. Number of ISTD Failures Allowed

Data Results:
Analytes: Pass
ISTD: Pass

Calibration Standard QC Report

Data File: C:\ICPCHEM\1\DATA\11K11100.B\006CALS.D\006CALS.D#
Date Acquired: Nov 11 2011 12:20 pm
Operator: NBS
Sample Name: 111111 Standard 2
Misc Info:
Vial Number: 1104
Current Method: C:\ICPCHEM\1\METHODS\62A1111A.M
Calibration File: C:\ICPCHEM\1\CALIB\62A1111A.C
Last Cal Update: Nov 11 2011 12:18 pm
Sample Type: CalStd
Total Dil Factor: 1.00

QC&ISTD Elements

Element	CPS Mean		SD	RSD(%)	Cal Coef
6 Li	3013436.00	A	14250.00	0.47	0.0000
7 (Li)	162843.41	P	655.10	0.40	1.0000
9 Be	10180.43	P	411.30	4.04	1.0000
11 B	16379.42	P	483.40	2.95	1.0000
23 Na	196689.50	P	7056.00	3.59	1.0000
24 Mg	23141.91	P	43.26	0.19	1.0000
27 Al	4021.80	P	226.90	5.64	1.0000
39 K	76357.12	P	2463.00	3.23	1.0000
44 Ca	1793.11	P	71.50	3.99	1.0000
45 Sc	510541.00	P	4569.00	0.89	0.0000
45 Sc	97262.66	P	635.50	0.65	0.0000
45 Sc	1465690.00	A	21530.00	1.47	0.0000
47 Ti	156.45	P	19.06	12.18	1.0000
51 V	8092.54	P	134.80	1.67	1.0000
52 Cr	4117.09	P	42.23	1.03	1.0000
55 Mn	61442.06	P	651.50	1.06	1.0000
56 Fe	82436.35	P	925.30	1.12	1.0000
59 Co	6109.79	P	52.36	0.86	1.0000
60 Ni	1383.64	P	28.30	2.05	1.0000
63 Cu	15516.40	P	233.60	1.51	1.0000
65 Cu	7559.83	P	73.09	0.97	1.0000
66 Zn	1430.31	P	74.87	5.23	1.0000
72 Ge	96818.69	P	1004.00	1.04	0.0000
72 Ge	44609.64	P	326.50	0.73	0.0000
72 Ge	203708.30	P	1751.00	0.86	0.0000
75 As	639.35	P	17.53	2.74	1.0000
78 Se	175.22	P	7.34	4.19	1.0000
78 Se	81.11	P	6.83	8.43	1.0000
88 Sr	3138.24	P	234.10	7.46	1.0000
88 Sr	17034.09	P	556.30	3.27	1.0000
95 Mo	3096.01	P	35.02	1.13	1.0000
106 (Cd)	180.01	P	18.56	10.31	1.0000
107 Ag	4028.49	P	77.05	1.91	1.0000
108 (Cd)	138.89	P	13.47	9.70	1.0000
111 Cd	1685.97	P	41.67	2.47	1.0000
115 In	1324038.00	A	6932.00	0.52	0.0000
118 Sn	5423.48	P	180.10	3.32	1.0000
121 Sb	6328.31	P	130.20	2.06	1.0000
137 Ba	2328.06	P	139.60	6.00	1.0000
159 Tb	1820559.00	A	17780.00	0.98	0.0000
165 Ho	1818461.00	A	19460.00	1.07	0.0000
205 Tl	15160.28	P	220.10	1.45	1.0000
206 (Pb)	7664.74	P	91.74	1.20	1.0000
207 (Pb)	7014.34	P	72.68	1.04	1.0000
208 Pb	31156.53	P	401.40	1.29	1.0000

ISTD Elements

Element	CPS Mean	RSD(%)	Ref Value	Rec(%)	QC Range(%)		Flag
6 Li	3013436.30	0.47	2775704.50	108.6	70 -	120	
45 Sc	510541.06	0.89	500780.41	101.9	70 -	120	
45 Sc	97262.66	0.65	95494.08	101.9	70 -	120	
45 Sc	1465690.00	1.47	1460980.80	100.3	70 -	120	
72 Ge	96818.70	1.04	96219.04	100.6	70 -	120	
72 Ge	44609.64	0.73	43611.78	102.3	70 -	120	
72 Ge	203708.33	0.86	213204.63	95.5	70 -	120	
115 In	1324038.00	0.52	1381264.00	95.9	70 -	120	
159 Tb	1820559.10	0.98	1843940.90	98.7	70 -	120	
165 Ho	1818460.60	1.07	1844184.90	98.6	70 -	120	

ISTD Ref File : C:\ICPCHEM\1\DATA\11K11100.B\004CALB.D\004CALB.D#

--- :Element Failures --- :Max. Number of Failures Allowed
0 :ISTD Failures 0 :Max. Number of ISTD Failures Allowed

Data Results:
Analytes: Pass
ISTD: Pass

Calibration Standard QC Report

Data File: C:\ICPCHEM\1\DATA\11K11100.B\007CALS.D\007CALS.D#
Date Acquired: Nov 11 2011 12:27 pm
Operator: NBS
Sample Name: 111111 Standard 3
Misc Info:
Vial Number: 1105
Current Method: C:\ICPCHEM\1\METHOD8\62A1111A.M
Calibration File: C:\ICPCHEM\1\CALIB\62A1111A.C
Last Cal Update: Nov 11 2011 12:24 pm
Sample Type: CalStd
Total Dil Factor: 1.00

QC&ISTD Elements

Element	CPS Mean		SD	RSD(%)	Cal Coef
6 Li	3129745.00	A	45250.00	1.45	0.0000
7 (Li)	169858.91	P	576.50	0.34	0.7236
9 Be	529831.50	P	2404.00	0.45	0.9999
11 B	352525.41	P	4097.00	1.16	0.9985
23 Na	1470286.00	A	17010.00	1.16	0.9979
24 Mg	1292682.00	A	17130.00	1.33	1.0000
27 Al	182025.00	P	2356.00	1.29	1.0000
39 K	502725.09	P	1262.00	0.25	0.9983
44 Ca	52265.13	P	789.60	1.51	0.9983
45 Sc	522323.31	P	4813.00	0.92	0.0000
45 Sc	98761.96	P	1402.00	1.42	0.0000
45 Sc	1523925.00	A	17440.00	1.14	0.0000
47 Ti	6316.10	P	62.52	0.99	0.9998
51 V	161320.00	P	2272.00	1.41	0.9994
52 Cr	179336.20	P	1262.00	0.70	1.0000
55 Mn	136966.41	P	806.40	0.59	0.9998
56 Fe	3466730.00	A	34680.00	1.00	1.0000
59 Co	254063.59	P	2599.00	1.02	0.9995
60 Ni	64869.85	P	659.40	1.02	0.9997
63 Cu	172209.41	P	983.40	0.57	0.9999
65 Cu	82567.48	P	346.90	0.42	1.0000
66 Zn	30294.80	P	353.70	1.17	0.9973
72 Ge	98255.19	P	550.50	0.56	0.0000
72 Ge	46262.58	P	34.38	0.07	0.0000
72 Ge	211131.20	P	2095.00	0.99	0.0000
75 As	20258.79	P	48.21	0.24	1.0000
78 Se	8196.34	P	137.70	1.68	1.0000
78 Se	2352.20	P	19.65	0.84	0.9963
88 Sr	152226.41	P	2676.00	1.76	0.9999
88 Sr	853159.19	P	3826.00	0.45	1.0000
95 Mo	152546.09	P	1308.00	0.86	0.9999
106 (Cd)	7779.08	P	43.36	0.56	0.9995
107 Ag	203275.00	P	1362.00	0.67	1.0000
108 (Cd)	5850.30	P	115.70	1.98	0.9966
111 Cd	85595.10	P	417.50	0.49	1.0000
115 In	1359449.00	A	15030.00	1.11	0.0000
118 Sn	233787.30	P	2145.00	0.92	0.9998
121 Sb	303264.81	P	1162.00	0.38	1.0000
137 Ba	112289.00	P	1153.00	1.03	1.0000
159 Tb	1852128.00	A	3859.00	0.21	0.0000
165 Ho	1866389.00	A	18420.00	0.99	0.0000
205 Tl	767163.63	P	3647.00	0.48	1.0000
206 (Pb)	267422.81	P	439.20	0.16	0.9998
207 (Pb)	229702.30	P	967.40	0.42	0.9996
208 Pb	1066559.00	P	3421.00	0.32	0.9997

ISTD Elements

Element	CPS Mean	RSD(%)	Ref Value	Rec(%)	QC Range(%)		Flag
6 Li	3129745.00	1.45	2775704.50	112.8	70 -	120	
45 Sc	522323.34	0.92	500780.41	104.3	70 -	120	
45 Sc	98761.96	1.42	95494.08	103.4	70 -	120	
45 Sc	1523925.40	1.14	1460980.80	104.3	70 -	120	
72 Ge	98255.19	0.56	96219.04	102.1	70 -	120	
72 Ge	46262.59	0.07	43611.78	106.1	70 -	120	
72 Ge	211131.19	0.99	213204.63	99.0	70 -	120	
115 In	1359449.00	1.11	1381264.00	98.4	70 -	120	
159 Tb	1852128.10	0.21	1843940.90	100.4	70 -	120	
165 Ho	1866389.00	0.99	1844184.90	101.2	70 -	120	

ISTD Ref File : C:\ICPCHEM\1\DATA\11K11100.B\004CALB.D\004CALB.D#

--- :Element Failures --- :Max. Number of Failures Allowed
0 :ISTD Failures 0 :Max. Number of ISTD Failures Allowed

Data Results:
Analytes: Pass
ISTD: Pass

Calibration Standard QC Report

Data File: C:\ICPCHEM\1\DATA\11K11100.B\008CALS.D\008CALS.D#
Date Acquired: Nov 11 2011 12:33 pm
Operator: NBS
Sample Name: 111111 Standard 4
Misc Info:
Vial Number: 1106
Current Method: C:\ICPCHEM\1\METHODS\62A1111A.M
Calibration File: C:\ICPCHEM\1\CALIB\62A1111A.C
Last Cal Update: Nov 11 2011 12:30 pm
Sample Type: CalStd
Total Dil Factor: 1.00

QC&ISTD Elements

Element	CPS Mean		SD	RSD(%)	Cal Coef
6 Li	3091825.00	A	34660.00	1.12	0.0000
7 (Li)	169207.09	P	2476.00	1.46	0.7966
9 Be	1184909.00	A	5168.00	0.44	1.0000
11 B	825041.63	A	8516.00	1.03	1.0000
23 Na	2686206.00	A	10730.00	0.40	0.9984
24 Mg	2535966.00	A	10300.00	0.41	1.0000
27 Al	366543.31	P	3283.00	0.90	1.0000
39 K	1039409.00	A	8793.00	0.85	0.9999
44 Ca	104136.20	P	1221.00	1.17	1.0000
45 Sc	526807.38	P	1501.00	0.28	0.0000
45 Sc	100637.20	P	272.50	0.27	0.0000
45 Sc	1546820.00	A	41280.00	2.67	0.0000
47 Ti	12883.09	P	335.40	2.60	1.0000
51 V	324487.59	P	1452.00	0.45	1.0000
52 Cr	360663.69	P	2389.00	0.66	1.0000
55 Mn	247566.30	P	2862.00	1.16	0.9063
56 Fe	6831163.00	A	89870.00	1.32	1.0000
59 Co	505973.59	P	1092.00	0.22	1.0000
60 Ni	128756.60	P	486.80	0.38	1.0000
63 Cu	331294.91	P	1236.00	0.37	0.9984
65 Cu	158678.41	P	595.70	0.38	0.9983
66 Zn	58476.31	P	247.90	0.42	0.9998
72 Ge	100101.50	P	582.20	0.58	0.0000
72 Ge	46752.66	P	94.61	0.20	0.0000
72 Ge	215920.09	P	4942.00	2.29	0.0000
75 As	41314.20	P	335.50	0.81	1.0000
78 Se	16782.86	P	111.00	0.66	1.0000
78 Se	4841.04	P	45.62	0.94	1.0000
88 Sr	308415.19	P	2179.00	0.71	1.0000
88 Sr	1836004.00	A	12020.00	0.65	1.0000
95 Mo	300376.41	P	620.60	0.20	1.0000
106 (Cd)	15606.90	P	85.03	0.54	1.0000
107 Ag	402429.69	P	2133.00	0.53	1.0000
108 (Cd)	11351.61	P	175.20	1.54	1.0000
111 Cd	169137.09	P	1111.00	0.66	1.0000
115 In	1356694.00	A	39030.00	2.88	0.0000
118 Sn	461432.81	P	1252.00	0.27	1.0000
121 Sb	616792.50	P	2811.00	0.46	1.0000
137 Ba	224905.80	P	424.60	0.19	1.0000
159 Tb	1896056.00	A	51090.00	2.69	0.0000
165 Ho	1892444.00	A	47210.00	2.49	0.0000
205 Tl	1621888.00	A	15450.00	0.95	1.0000
206 (Pb)	524239.50	P	2392.00	0.46	1.0000
207 (Pb)	454785.81	P	2844.00	0.63	1.0000
208 Pb	2164409.00	A	4337.00	0.20	1.0000

ISTD Elements

Element	CPS Mean	RSD(%)	Ref Value	Rec(%)	QC Range(%)		Flag
6 Li	3091824.50	1.12	2775704.50	111.4	70 -	120	
45 Sc	526807.44	0.28	500780.41	105.2	70 -	120	
45 Sc	100637.22	0.27	95494.08	105.4	70 -	120	
45 Sc	1546819.60	2.67	1460980.80	105.9	70 -	120	
72 Ge	100101.52	0.58	96219.04	104.0	70 -	120	
72 Ge	46752.66	0.20	43611.78	107.2	70 -	120	
72 Ge	215920.11	2.29	213204.63	101.3	70 -	120	
115 In	1356693.50	2.88	1381264.00	98.2	70 -	120	
159 Tb	1896055.90	2.69	1843940.90	102.8	70 -	120	
165 Ho	1892443.90	2.49	1844184.90	102.6	70 -	120	

ISTD Ref File : C:\ICPCHEM\1\DATA\11K11100.B\004CALB.D\004CALB.D#

--- :Element Failures --- :Max. Number of Failures Allowed
0 :ISTD Failures 0 :Max. Number of ISTD Failures Allowed

Data Results:
Analytes: Pass
ISTD: Pass

QCS QC Report

Data File: C:\ICPCHEM\1\DATA\11K11100.B\009_QCS.D\009_QCS.D#
Date Acquired: Nov 11 2011 12:39 pm
Operator: NBS
Sample Name: ICV 111111
Misc Info:
Vial Number: 1107
Current Method: C:\ICPCHEM\1\METHODS\62A1111A.M
Calibration File: C:\ICPCHEM\1\CALIB\62A1111A.C
Last Cal Update: Nov 11 2011 12:36 pm
Sample Type: QCS
Total Dil Factor: 1.00

QC Elements

Element	Conc.		RSD(%)	Expected	QC Range(%)		Flag
7 (Li)	--------	ug/l	--------	100.00	90 -	110	
9 Be	107.60	ug/l	0.95	100.00	90 -	110	
11 B	105.70	ug/l	0.67	100.00	90 -	110	
23 Na	2518.00	ug/l	0.81	2500.00	90 -	110	
24 Mg	2533.00	ug/l	0.67	2500.00	90 -	110	
27 Al	2547.00	ug/l	1.32	2500.00	90 -	110	
39 K	2615.00	ug/l	0.71	2500.00	90 -	110	
44 Ca	2519.00	ug/l	0.47	2500.00	90 -	110	
47 Ti	97.29	ug/l	0.90	100.00	90 -	110	
51 V	103.40	ug/l	0.55	100.00	90 -	110	
52 Cr	106.50	ug/l	0.63	100.00	90 -	110	
55 Mn	106.70	ug/l	0.21	100.00	90 -	110	
56 Fe	2516.00	ug/l	1.06	2500.00	90 -	110	
59 Co	104.60	ug/l	0.25	100.00	90 -	110	
60 Ni	104.70	ug/l	0.28	100.00	90 -	110	
63 Cu	102.50	ug/l	1.70	100.00	90 -	110	
65 Cu	102.20	ug/l	1.45	100.00	90 -	110	
66 Zn	104.10	ug/l	1.10	100.00	90 -	110	
75 As	98.86	ug/l	1.38	100.00	90 -	110	
78 Se	103.60	ug/l	1.81	100.00	90 -	110	
78 Se	104.10	ug/l	2.03	100.00	90 -	110	
88 Sr	101.20	ug/l	1.63	100.00	90 -	110	
88 Sr	104.30	ug/l	0.60	100.00	90 -	110	
95 Mo	96.15	ug/l	1.35	100.00	90 -	110	
106 (Cd)	--------	ug/l	--------	100.00	90 -	110	
107 Ag	46.26	ug/l	0.71	50.00	90 -	110	
108 (Cd)	--------	ug/l	--------	100.00	90 -	110	
111 Cd	103.60	ug/l	0.47	100.00	90 -	110	
118 Sn	43.82	ug/l	0.17	50.00	90 -	110	Fail
121 Sb	102.70	ug/l	0.18	100.00	90 -	110	
137 Ba	99.56	ug/l	0.34	100.00	90 -	110	
205 Tl	106.40	ug/l	1.20	100.00	90 -	110	
206 (Pb)	--------	ug/l	--------	100.00	90 -	110	
207 (Pb)	--------	ug/l	--------	100.00	90 -	110	
208 Pb	106.30	ug/l	0.89	100.00	90 -	110	

ISTD Elements

Element	CPS Mean	RSD(%)	Ref Value	Rec(%)	QC Range(%)		Flag
6 Li	3157481.00	0.39	2775704.50	113.8	70 -	120	
45 Sc	523431.13	0.16	500780.41	104.5	70 -	120	
45 Sc	100384.52	0.40	95494.08	105.1	70 -	120	
45 Sc	1532510.60	0.50	1460980.80	104.9	70 -	120	
72 Ge	99727.78	0.25	96219.04	103.6	70 -	120	
72 Ge	46938.75	0.91	43611.78	107.6	70 -	120	
72 Ge	212937.78	0.32	213204.63	99.9	70 -	120	
115 In	1371120.50	0.09	1381264.00	99.3	70 -	120	
159 Tb	1873353.00	0.83	1843940.90	101.6	70 -	120	
165 Ho	1868336.50	1.05	1844184.90	101.3	70 -	120	

ISTD Ref File : C:\ICPCHEM\1\DATA\11K11100.B\004CALB.D\004CALB.D#

1 :Element Failures 0 :Max. Number of Failures Allowed
0 :ISTD Failures 0 :Max. Number of ISTD Failures Allowed

Data Results:
Analytes: Fail
ISTD: Pass

CCB QC Report

Data File: C:\ICPCHEM\1\DATA\11K11100.B\012_CCB.D\012_CCB.D#
Date Acquired: Nov 11 2011 12:57 pm
Operator: NBS
Sample Name: ICB 111111
Misc Info:
Vial Number: 1102
Current Method: C:\ICPCHEM\1\METHODS\62A1111A.M
Calibration File: C:\ICPCHEM\1\CALIB\62A1111A.C
Last Cal Update: Nov 11 2011 12:36 pm
Sample Type: CCB
Total Dil Factor: 1.00

QC Elements

Element	Conc.	RSD(%)	High Limit	Flag
7 (Li)	-------- ug/l	--------	#VALUE!	
9 Be	0.00 ug/l	64.25	0.12	
11 B	0.03 ug/l	45.15	15.00	
23 Na	7.71 ug/l	8.18	77.10	
24 Mg	0.10 ug/l	55.44	7.50	
27 Al	0.09 ug/l	51.89	3.96	
39 K	-16.07 ug/l	31.35	19.20	
44 Ca	2.26 ug/l	102.26	90.00	
47 Ti	0.02 ug/l	221.48	0.78	
51 V	0.57 ug/l	2.59	0.21	Fail
52 Cr	0.01 ug/l	92.18	0.12	
55 Mn	0.00 ug/l	249.24	0.18	
56 Fe	0.25 ug/l	4.89	40.80	
59 Co	-0.25 ug/l	1.58	0.09	
60 Ni	0.00 ug/l	280.61	0.48	
63 Cu	-0.13 ug/l	3.22	0.39	
65 Cu	-0.13 ug/l	16.74	0.39	
66 Zn	-0.01 ug/l	406.21	6.90	
75 As	-0.09 ug/l	15.13	0.27	
78 Se	0.01 ug/l	58.10	0.30	
78 Se	0.05 ug/l	139.53	0.30	
88 Sr	0.00 ug/l	1034.40	0.03	
88 Sr	0.00 ug/l	24.09	0.03	
95 Mo	0.03 ug/l	16.72	0.21	
106 (Cd)	-------- ug/l	--------	#VALUE!	
107 Ag	0.00 ug/l	50.39	0.09	
108 (Cd)	-------- ug/l	--------	#VALUE!	
111 Cd	0.01 ug/l	58.78	0.06	
118 Sn	0.03 ug/l	55.51	0.30	
121 Sb	0.13 ug/l	5.57	0.03	Fail
137 Ba	0.01 ug/l	116.79	0.12	
205 Tl	0.01 ug/l	38.28	0.03	
206 (Pb)	-------- ug/l	--------	#VALUE!	
207 (Pb)	-------- ug/l	--------	#VALUE!	
208 Pb	-0.20 ug/l	0.78	0.33	

ISTD Elements

Element	CPS Mean	RSD(%)	Ref Value	Rec(%)	QC Range(%)	Flag
6 Li	3073279.00	0.84	2775704.50	110.7	70 - 120	
45 Sc	545909.38	3.12	500780.41	109.0	70 - 120	
45 Sc	100165.70	0.44	95494.08	104.9	70 - 120	
45 Sc	1499557.30	0.22	1460980.80	102.6	70 - 120	
72 Ge	101795.60	2.62	96219.04	105.8	70 - 120	
72 Ge	46734.16	0.18	43611.78	107.2	70 - 120	
72 Ge	210654.83	0.54	213204.63	98.8	70 - 120	
115 In	1336860.30	0.89	1381264.00	96.8	70 - 120	
159 Tb	1857728.00	1.11	1843940.90	100.7	70 - 120	
165 Ho	1856236.60	1.27	1844184.90	100.7	70 - 120	

ISTD Ref File : C:\ICPCHEM\1\DATA\11K11100.B\004CALB.D\004CALB.D#

2 :Element Failures 0 :Max. Number of Failures Allowed
0 :ISTD Failures 0 :Max. Number of ISTD Failures Allowed

Data Results:
Analytes: Fail
ISTD: Pass

CCV QC Report

Data File: C:\ICPCHEM\1\DATA\11K11100.B\013_CCV.D\013_CCV.D#
Date Acquired: Nov 11 2011 01:03 pm
Operator: NBS
Sample Name: CCV 111111
Misc Info:
Vial Number: 1105
Current Method: C:\ICPCHEM\1\METHODS\62A1111A.M
Calibration File: C:\ICPCHEM\1\CALIB\62A1111A.C
Last Cal Update: Nov 11 2011 12:36 pm
Sample Type: CCV
Total Dil Factor: 1.00

QC Elements

Element	Conc.		RSD(%)	Expected	QC Range(%)		Flag
7 (Li)	--------	ug/l	--------	50.00	90 -	110	
9 Be	45.86	ug/l	1.99	50.00	90 -	110	
11 B	43.69	ug/l	2.15	50.00	90 -	110	Fail
23 Na	1276.00	ug/l	1.53	1250.00	90 -	110	
24 Mg	2559.00	ug/l	1.05	2500.00	90 -	110	
27 Al	1001.00	ug/l	1.70	1000.00	90 -	110	
39 K	917.90	ug/l	1.26	1000.00	90 -	110	
44 Ca	2498.00	ug/l	1.41	2500.00	90 -	110	
47 Ti	49.40	ug/l	0.95	50.00	90 -	110	
51 V	50.61	ug/l	1.18	50.00	90 -	110	
52 Cr	50.27	ug/l	1.38	50.00	90 -	110	
55 Mn	54.78	ug/l	1.56	50.00	90 -	110	
56 Fe	1027.00	ug/l	1.77	1000.00	90 -	110	
59 Co	50.74	ug/l	0.93	50.00	90 -	110	
60 Ni	50.88	ug/l	1.81	50.00	90 -	110	
63 Cu	50.81	ug/l	0.58	50.00	90 -	110	
65 Cu	50.69	ug/l	0.50	50.00	90 -	110	
66 Zn	51.32	ug/l	0.03	50.00	90 -	110	
75 As	49.12	ug/l	0.62	50.00	90 -	110	
78 Se	50.32	ug/l	1.71	50.00	90 -	110	
78 Se	49.06	ug/l	1.10	50.00	90 -	110	
88 Sr	49.93	ug/l	0.31	50.00	90 -	110	
88 Sr	47.83	ug/l	0.44	50.00	90 -	110	
95 Mo	50.08	ug/l	0.70	50.00	90 -	110	
106 (Cd)	--------	ug/l	--------	50.00	90 -	110	
107 Ag	24.63	ug/l	0.57	25.00	90 -	110	
108 (Cd)	--------	ug/l	--------	50.00	90 -	110	
111 Cd	50.07	ug/l	1.49	50.00	90 -	110	
118 Sn	50.35	ug/l	0.28	50.00	90 -	110	
121 Sb	49.48	ug/l	0.77	50.00	90 -	110	
137 Ba	49.18	ug/l	0.66	50.00	90 -	110	
205 Tl	48.89	ug/l	0.40	50.00	90 -	110	
206 (Pb)	--------	ug/l	--------	50.00	90 -	110	
207 (Pb)	--------	ug/l	--------	50.00	90 -	110	
208 Pb	50.31	ug/l	0.14	50.00	90 -	110	

ISTD Elements

Element	CPS Mean	RSD(%)	Ref Value	Rec(%)	QC Range(%)		Flag
6 Li	3009330.80	0.85	2775704.50	108.4	70 -	120	
45 Sc	502422.56	3.92	500780.41	100.3	70 -	120	
45 Sc	98428.88	1.42	95494.08	103.1	70 -	120	
45 Sc	1480640.80	1.02	1460980.80	101.3	70 -	120	
72 Ge	97237.93	2.52	96219.04	101.1	70 -	120	
72 Ge	46537.16	0.17	43611.78	106.7	70 -	120	
72 Ge	206334.70	0.20	213204.63	96.8	70 -	120	
115 In	1333758.10	0.36	1381264.00	96.6	70 -	120	
159 Tb	1832635.60	0.51	1843940.90	99.4	70 -	120	
165 Ho	1824652.90	0.58	1844184.90	98.9	70 -	120	

ISTD Ref File : C:\ICPCHEM\1\DATA\11K11100.B\004CALB.D\004CALB.D#

1 :Element Failures 0 :Max. Number of Failures Allowed
0 :ISTD Failures 0 :Max. Number of ISTD Failures Allowed

Data Results:
Analytes: Fail
ISTD: Pass

CCB QC Report

Data File: C:\ICPCHEM\1\DATA\11K11100.B\014_CCB.D\014_CCB.D#
Date Acquired: Nov 11 2011 01:09 pm
Operator: NBS
Sample Name: CCB 111111
Misc Info:
Vial Number: 1102
Current Method: C:\ICPCHEM\1\METHODS\62A1111A.M
Calibration File: C:\ICPCHEM\1\CALIB\62A1111A.C
Last Cal Update: Nov 11 2011 12:36 pm
Sample Type: CCB
Total Dil Factor: 1.00

QC Elements

Element	Conc.	RSD(%)	High Limit	Flag
7 (Li)	-------- ug/l	--------	#VALUE!	
9 Be	0.00 ug/l	75.05	0.12	
11 B	0.19 ug/l	2.24	15.00	
23 Na	3.09 ug/l	28.12	77.10	
24 Mg	0.17 ug/l	59.53	7.50	
27 Al	0.11 ug/l	49.92	3.96	
39 K	-13.77 ug/l	41.47	19.20	
44 Ca	1.49 ug/l	82.56	90.00	
47 Ti	0.00 ug/l	584.34	0.78	
51 V	0.76 ug/l	2.01	0.21	Fail
52 Cr	0.02 ug/l	22.00	0.12	
55 Mn	0.01 ug/l	26.62	0.18	
56 Fe	0.44 ug/l	10.30	40.80	
59 Co	-0.27 ug/l	0.49	0.09	
60 Ni	0.00 ug/l	169.21	0.48	
63 Cu	-0.16 ug/l	12.22	0.39	
65 Cu	-0.16 ug/l	4.20	0.39	
66 Zn	0.03 ug/l	67.54	6.90	
75 As	-0.03 ug/l	60.04	0.27	
78 Se	0.10 ug/l	23.77	0.30	
78 Se	0.02 ug/l	138.81	0.30	
88 Sr	0.00 ug/l	574.89	0.03	
88 Sr	0.00 ug/l	33.45	0.03	
95 Mo	0.10 ug/l	2.35	0.21	
106 (Cd)	-------- ug/l	--------	#VALUE!	
107 Ag	0.00 ug/l	6.78	0.09	
108 (Cd)	-------- ug/l	--------	#VALUE!	
111 Cd	0.01 ug/l	137.32	0.06	
118 Sn	0.06 ug/l	31.67	0.30	
121 Sb	0.69 ug/l	6.56	0.03	Fail
137 Ba	0.01 ug/l	127.23	0.12	
205 Tl	0.02 ug/l	7.09	0.03	
206 (Pb)	-------- ug/l	--------	#VALUE!	
207 (Pb)	-------- ug/l	--------	#VALUE!	
208 Pb	-0.21 ug/l	0.72	0.33	

ISTD Elements

Element	CPS Mean	RSD(%)	Ref Value	Rec(%)	QC Range(%)	Flag
6 Li	2980962.00	0.72	2775704.50	107.4	70 - 120	
45 Sc	505025.28	3.19	500780.41	100.8	70 - 120	
45 Sc	97675.82	0.75	95494.08	102.3	70 - 120	
45 Sc	1485366.30	0.36	1460980.80	101.7	70 - 120	
72 Ge	97202.46	2.05	96219.04	101.0	70 - 120	
72 Ge	45665.85	0.21	43611.78	104.7	70 - 120	
72 Ge	205716.23	0.30	213204.63	96.5	70 - 120	
115 In	1321174.40	0.50	1381264.00	95.6	70 - 120	
159 Tb	1807747.90	0.40	1843940.90	98.0	70 - 120	
165 Ho	1813776.00	0.73	1844184.90	98.4	70 - 120	

ISTD Ref File : C:\ICPCHEM\1\DATA\11K11100.B\004CALB.D\004CALB.D#

2 :Element Failures 0 :Max. Number of Failures Allowed
0 :ISTD Failures 0 :Max. Number of ISTD Failures Allowed

Data Results:
Analytes: Fail
ISTD: Pass

ICS-A QC Report

Data File: C:\ICPCHEM\1\DATA\11K11100.B\015ICSA.D\015ICSA.D#
Date Acquired: Nov 11 2011 01:15 pm
Acq. Method: 62A1111A.M
Operator: NBS
Sample Name: ICSA 111111
Misc Info:
Vial Number: 2102
Current Method: C:\ICPCHEM\1\METHODS\62A1111A.M
Calibration File: C:\ICPCHEM\1\CALIB\62A1111A.C
Last Cal. Update: Nov 11 2011 12:36 pm
Sample Type: ICSA
Dilution Factor: 1.00

Data Results:
Analytes: Pass
ISTD: Pass

QC Elements

Element	IS Ref	Tune	Conc.		RSD(%)	High Limit ppb	Flag
7 (Li)	---	3	--------	ug/l	--------		
9 Be	45	3	0.83	ug/l	2.55		
11 B	45	3	1.56	ug/l	2.76		
23 Na	45	2	93250.00	ug/l	0.74		
24 Mg	45	2	91170.00	ug/l	1.03		
27 Al	45	2	104800.00	ug/l	1.31		
39 K	45	2	95010.00	ug/l	0.86		
44 Ca	45	2	101900.00	ug/l	1.03		
47 Ti	45	2	1961.00	ug/l	0.73		
51 V	45	2	2.53	ug/l	1.39		
52 Cr	45	2	2.35	ug/l	1.68		
55 Mn	45	2	7.50	ug/l	0.84		
56 Fe	45	2	92610.00	ug/l	0.44		
59 Co	45	2	20.49	ug/l	0.58		
60 Ni	45	2	3.86	ug/l	0.68		
63 Cu	72	2	1.60	ug/l	2.77		
65 Cu	72	2	1.70	ug/l	4.01		
66 Zn	72	2	5.11	ug/l	1.25		
75 As	72	2	1.55	ug/l	3.09		
78 Se	72	1	1.07	ug/l	5.94		
78 Se	72	2	1.16	ug/l	9.00		
88 Sr	72	2	1.41	ug/l	4.62		
88 Sr	72	3	1.37	ug/l	1.32		
95 Mo	72	3	1834.00	ug/l	1.74		
106 (Cd)	---	3	--------	ug/l	--------		
107 Ag	115	3	1.97	ug/l	1.05		
108 (Cd)	---	3	--------	ug/l	--------		
111 Cd	115	3	2.42	ug/l	3.67		
118 Sn	115	3	1.18	ug/l	1.50		
121 Sb	115	3	1.93	ug/l	2.34		
137 Ba	115	3	3.88	ug/l	2.04		
205 Tl	159	3	1.62	ug/l	1.90		
206 (Pb)	---	3	--------	ug/l	--------		
207 (Pb)	---	3	--------	ug/l	--------		
208 Pb	159	3	3.50	ug/l	0.41		

ISTD Elements

Element	Tune	CPS Mean	RSD(%)	Ref Value	Rec(%)	QC Range(%)	Flag
6 Li	3	2777926	0.73	2775705	100.1	70 - 120	
45 Sc	1	527513	3.31	500780	105.3	70 - 120	
45 Sc	2	94664	0.69	95494	99.1	70 - 120	
45 Sc	3	1465735	0.50	1460981	100.3	70 - 120	
72 Ge	1	98457	2.56	96219	102.3	70 - 120	
72 Ge	2	46798	1.22	43612	107.3	70 - 120	
72 Ge	3	216093	0.53	213205	101.4	70 - 120	
115 In	3	1235992	0.56	1381264	89.5	70 - 120	
159 Tb	3	1778881	0.42	1843941	96.5	70 - 120	
165 Ho	3	1783575	1.04	1844185	96.7	70 - 120	

Tune File# 1 c:\icpchem\1\7500\h2.u
Tune File# 2 c:\icpchem\1\7500\he.u
Tune File# 3 c:\icpchem\1\7500\nogas.u

ISTD Ref File : C:\ICPCHEM\1\DATA\11K11100.B\004CALB.D\004CALB.D#

0 :Element Failures
0 :ISTD Failures
0 :Max. Number of Failures Allowed
0 :Max. Nnumber of ISTD Failures Allowed

ICS-AB QC Report

Data File: C:\ICPCHEM\1\DATA\11K11100.B\016ICSB.D\016ICSB.D#
Date Acquired: Nov 11 2011 01:21 pm
Acq. Method: 62A1111A.M
Operator: NBS
Sample Name: ICSAB 111111
Misc Info:
Vial Number: 2103
Current Method: C:\ICPCHEM\1\METHODS\62A1111A.M
Calibration File: C:\ICPCHEM\1\CALIB\62A1111A.C
Last Cal. Update: Nov 11 2011 12:36 pm
Sample Type: ICSAB
Dilution Factor: 1.00

Data Results:
Analytes: Pass
ISTD: Pass

QC Elements

Element	IS Ref	Tune	Conc. ppb	RSD(%)	Expected	%Recovery	QC Range(%)	Flag
7 (Li)	---	3	--------	--------	---		-	
9 Be	45	3	246.60	1.11	250	98.6	80 - 120	
11 B	45	3	1.63	0.34	---		-	
23 Na	45	2	96100.00	0.19	---		-	
24 Mg	45	2	93890.00	0.39	---		-	
27 Al	45	2	107500.00	1.06	---		-	
39 K	45	2	97710.00	0.66	---		-	
44 Ca	45	2	105500.00	1.05	---		-	
47 Ti	45	2	2014.00	0.31	2000	100.7	80 - 120	
51 V	45	2	267.30	0.98	250	106.9	80 - 120	
52 Cr	45	2	270.40	2.94	250	108.2	80 - 120	
55 Mn	45	2	264.30	0.56	250	105.7	80 - 120	
56 Fe	45	2	94360.00	0.20	---		-	
59 Co	45	2	282.40	0.78	250	113.0	80 - 120	
60 Ni	45	2	481.90	0.90	500	96.4	80 - 120	
63 Cu	72	2	218.20	1.17	250	87.3	80 - 120	
65 Cu	72	2	218.60	0.91	250	87.4	80 - 120	
66 Zn	72	2	513.60	0.37	500	102.7	80 - 120	
75 As	72	2	239.20	0.52	250	95.7	80 - 120	
78 Se	72	1	251.50	0.86	250	100.6	80 - 120	
78 Se	72	2	233.50	0.80	250	93.4	80 - 120	
88 Sr	72	2	1.62	0.60	---		-	
88 Sr	72	3	1.51	0.87	---		-	
95 Mo	72	3	2131.00	0.52	2000	106.6	80 - 120	
106 (Cd)	---	3	--------	--------	---		-	
107 Ag	115	3	535.90	1.10	500	107.2	80 - 120	
108 (Cd)	---	3	--------	--------	---		-	
111 Cd	115	3	495.00	0.98	500	99.0	80 - 120	
118 Sn	115	3	1.45	2.79	---		-	
121 Sb	115	3	274.60	0.82	250	109.8	80 - 120	
137 Ba	115	3	271.00	0.98	250	108.4	80 - 120	
205 Tl	159	3	252.50	0.03	250	101.0	80 - 120	
206 (Pb)	---	3	--------	--------	---		-	
207 (Pb)	---	3	--------	--------	---		-	
208 Pb	159	3	502.00	0.13	500	100.4	80 - 120	

ISTD Elements

Element	Tune	CPS Mean	RSD(%)	Ref Value	Rec(%)	QC Range(%)	Flag
6 Li	3	2732517	0.30	2775705	98.4	70 - 120	
45 Sc	1	511020	0.99	500780	102.0	70 - 120	
45 Sc	2	93932	0.16	95494	98.4	70 - 120	
45 Sc	3	1418244	0.91	1460981	97.1	70 - 120	
72 Ge	1	96432	0.78	96219	100.2	70 - 120	
72 Ge	2	46185	0.98	43612	105.9	70 - 120	
72 Ge	3	209601	0.65	213205	98.3	70 - 120	
115 In	3	1203221	0.93	1381264	87.1	70 - 120	
159 Tb	3	1775149	0.42	1843941	96.3	70 - 120	
165 Ho	3	1779108	0.39	1844185	96.5	70 - 120	

Tune File# 1 c:\icpchem\1\7500\h2.u
Tune File# 2 c:\icpchem\1\7500\he.u
Tune File# 3 c:\icpchem\1\7500\nogas.u

ISTD Ref File : C:\ICPCHEM\1\DATA\11K11100.B\004CALB.D\004CALB.D#

0 :Element Failures
0 :ISTD Failures
0 :Max. Number of Failures Allowed
0 :Max. Number of ISTD Failures Allowed

CCV QC Report

Data File: C:\ICPCHEM\1\DATA\11K11100.B\018_CCV.D\018_CCV.D#
Date Acquired: Nov 11 2011 01:33 pm
Operator: NBS
Sample Name: CCV 111111
Misc Info:
Vial Number: 1105
Current Method: C:\ICPCHEM\1\METHODS\62A1111A.M
Calibration File: C:\ICPCHEM\1\CALIB\62A1111A.C
Last Cal Update: Nov 11 2011 12:36 pm
Sample Type: CCV
Total Dil Factor: 1.00

QC Elements

Element	Conc.		RSD(%)	Expected	QC Range(%)		Flag
7 (Li)	--------	ug/l	--------	50.00	90 -	110	
9 Be	45.54	ug/l	0.43	50.00	90 -	110	
11 B	42.52	ug/l	0.69	50.00	90 -	110	Fail
23 Na	1232.00	ug/l	0.90	1250.00	90 -	110	
24 Mg	2553.00	ug/l	1.13	2500.00	90 -	110	
27 Al	1002.00	ug/l	1.24	1000.00	90 -	110	
39 K	910.10	ug/l	0.77	1000.00	90 -	110	
44 Ca	2506.00	ug/l	0.99	2500.00	90 -	110	
47 Ti	49.11	ug/l	2.38	50.00	90 -	110	
51 V	50.62	ug/l	1.42	50.00	90 -	110	
52 Cr	49.48	ug/l	1.20	50.00	90 -	110	
55 Mn	54.25	ug/l	1.47	50.00	90 -	110	
56 Fe	1014.00	ug/l	0.73	1000.00	90 -	110	
59 Co	50.51	ug/l	0.59	50.00	90 -	110	
60 Ni	50.70	ug/l	1.36	50.00	90 -	110	
63 Cu	49.70	ug/l	0.40	50.00	90 -	110	
65 Cu	49.73	ug/l	0.19	50.00	90 -	110	
66 Zn	49.96	ug/l	0.97	50.00	90 -	110	
75 As	48.64	ug/l	0.06	50.00	90 -	110	
78 Se	48.54	ug/l	0.17	50.00	90 -	110	
78 Se	48.82	ug/l	0.30	50.00	90 -	110	
88 Sr	49.77	ug/l	0.41	50.00	90 -	110	
88 Sr	48.29	ug/l	0.56	50.00	90 -	110	
95 Mo	51.00	ug/l	1.55	50.00	90 -	110	
106 (Cd)	--------	ug/l	--------	50.00	90 -	110	
107 Ag	25.03	ug/l	0.97	25.00	90 -	110	
108 (Cd)	--------	ug/l	--------	50.00	90 -	110	
111 Cd	49.94	ug/l	1.03	50.00	90 -	110	
118 Sn	51.03	ug/l	1.26	50.00	90 -	110	
121 Sb	50.47	ug/l	0.55	50.00	90 -	110	
137 Ba	50.01	ug/l	0.80	50.00	90 -	110	
205 Tl	48.64	ug/l	0.31	50.00	90 -	110	
206 (Pb)	--------	ug/l	--------	50.00	90 -	110	
207 (Pb)	--------	ug/l	--------	50.00	90 -	110	
208 Pb	50.34	ug/l	1.15	50.00	90 -	110	

ISTD Elements

Element	CPS Mean	RSD(%)	Ref Value	Rec(%)	QC Range(%)		Flag
6 Li	3132762.00	0.73	2775704.50	112.9	70 -	120	
45 Sc	534639.31	0.21	500780.41	106.8	70 -	120	
45 Sc	99077.84	0.49	95494.08	103.8	70 -	120	
45 Sc	1513729.60	0.46	1460980.80	103.6	70 -	120	
72 Ge	102211.11	0.47	96219.04	106.2	70 -	120	
72 Ge	47244.57	0.27	43611.78	108.3	70 -	120	
72 Ge	214737.88	0.60	213204.63	100.7	70 -	120	
115 In	1389034.00	1.09	1381264.00	100.6	70 -	120	
159 Tb	1908915.90	0.41	1843940.90	103.5	70 -	120	
165 Ho	1921136.40	0.48	1844184.90	104.2	70 -	120	

ISTD Ref File : C:\ICPCHEM\1\DATA\11K11100.B\004CALB.D\004CALB.D#

1 :Element Failures 0 :Max. Number of Failures Allowed
0 :ISTD Failures 0 :Max. Number of ISTD Failures Allowed

Data Results:
Analytes: Fail
ISTD: Pass

CCB QC Report

Data File: C:\ICPCHEM\1\DATA\11K11100.B\020_CCB.D\020_CCB.D#
Date Acquired: Nov 11 2011 01:46 pm
Operator: NBS
Sample Name: CCB 111111
Misc Info:
Vial Number: 1102
Current Method: C:\ICPCHEM\1\METHODS\62A1111A.M
Calibration File: C:\ICPCHEM\1\CALIB\62A1111A.C
Last Cal Update: Nov 11 2011 12:36 pm
Sample Type: CCB
Total Dil Factor: 1.00

QC Elements

Element	Conc.	RSD(%)	High Limit	Flag
7 (Li)	-------- ug/l	--------	#VALUE!	
9 Be	0.00 ug/l	342.89	0.12	
11 B	-0.21 ug/l	27.87	15.00	
23 Na	-22.55 ug/l	2.00	77.10	
24 Mg	0.32 ug/l	24.61	7.50	
27 Al	0.31 ug/l	30.94	3.96	
39 K	-13.04 ug/l	28.54	19.20	
44 Ca	-2.47 ug/l	54.66	90.00	
47 Ti	0.02 ug/l	115.52	0.78	
51 V	1.17 ug/l	4.59	0.21	Fail
52 Cr	0.01 ug/l	119.76	0.12	
55 Mn	0.00 ug/l	605.19	0.18	
56 Fe	0.83 ug/l	6.77	40.80	
59 Co	-0.30 ug/l	0.59	0.09	
60 Ni	0.00 ug/l	211.25	0.48	
63 Cu	-0.31 ug/l	1.69	0.39	
65 Cu	-0.31 ug/l	4.87	0.39	
66 Zn	0.00 ug/l	866.90	6.90	
75 As	-0.07 ug/l	27.14	0.27	
78 Se	0.03 ug/l	41.10	0.30	
78 Se	0.07 ug/l	100.62	0.30	
88 Sr	0.00 ug/l	114.02	0.03	
88 Sr	0.00 ug/l	18.44	0.03	
95 Mo	0.09 ug/l	4.01	0.21	
106 (Cd)	-------- ug/l	--------	#VALUE!	
107 Ag	0.01 ug/l	12.14	0.09	
108 (Cd)	-------- ug/l	--------	#VALUE!	
111 Cd	0.00 ug/l	1077.10	0.06	
118 Sn	0.03 ug/l	30.18	0.30	
121 Sb	0.29 ug/l	4.29	0.03	Fail
137 Ba	0.01 ug/l	115.49	0.12	
205 Tl	0.01 ug/l	19.13	0.03	
206 (Pb)	-------- ug/l	--------	#VALUE!	
207 (Pb)	-------- ug/l	--------	#VALUE!	
208 Pb	-0.21 ug/l	1.63	0.33	

ISTD Elements

Element	CPS Mean	RSD(%)	Ref Value	Rec(%)	QC Range(%)	Flag
6 Li	3042523.00	0.11	2775704.50	109.6	70 - 120	
45 Sc	529492.56	0.66	500780.41	105.7	70 - 120	
45 Sc	97690.27	0.93	95494.08	102.3	70 - 120	
45 Sc	1482243.40	0.75	1460980.80	101.5	70 - 120	
72 Ge	101254.01	0.60	96219.04	105.2	70 - 120	
72 Ge	46065.66	0.31	43611.78	105.6	70 - 120	
72 Ge	210454.86	0.84	213204.63	98.7	70 - 120	
115 In	1353362.30	0.71	1381264.00	98.0	70 - 120	
159 Tb	1859786.10	0.52	1843940.90	100.9	70 - 120	
165 Ho	1863063.90	0.81	1844184.90	101.0	70 - 120	

ISTD Ref File : C:\ICPCHEM\1\DATA\11K11100.B\004CALB.D\004CALB.D#

2 :Element Failures 0 :Max. Number of Failures Allowed
0 :ISTD Failures 0 :Max. Number of ISTD Failures Allowed

Data Results:
Analytes: Fail
ISTD: Pass

CCV QC Report

Data File: C:\ICPCHEM\1\DATA\11K11100.B\032_CCV.D\032_CCV.D#
Date Acquired: Nov 11 2011 03:05 pm
Operator: NBS
Sample Name: CCV 111111
Misc Info:
Vial Number: 1105
Current Method: C:\ICPCHEM\1\METHODS\62A1111A.M
Calibration File: C:\ICPCHEM\1\CALIB\62A1111A.C
Last Cal Update: Nov 11 2011 12:36 pm
Sample Type: CCV
Total Dil Factor: 1.00

QC Elements

Element	Conc.		RSD(%)	Expected	QC Range(%)		Flag
7 (Li)	--------	ug/l	--------	50.00	90 -	110	
9 Be	45.98	ug/l	0.90	50.00	90 -	110	
11 B	43.38	ug/l	1.83	50.00	90 -	110	Fail
23 Na	1250.00	ug/l	2.32	1250.00	90 -	110	
24 Mg	2541.00	ug/l	1.13	2500.00	90 -	110	
27 Al	992.10	ug/l	1.33	1000.00	90 -	110	
39 K	905.50	ug/l	1.76	1000.00	90 -	110	
44 Ca	2473.00	ug/l	1.61	2500.00	90 -	110	
47 Ti	49.01	ug/l	0.71	50.00	90 -	110	
51 V	50.64	ug/l	0.61	50.00	90 -	110	
52 Cr	49.61	ug/l	0.94	50.00	90 -	110	
55 Mn	54.01	ug/l	0.87	50.00	90 -	110	
56 Fe	1013.00	ug/l	1.52	1000.00	90 -	110	
59 Co	50.36	ug/l	0.94	50.00	90 -	110	
60 Ni	51.02	ug/l	1.41	50.00	90 -	110	
63 Cu	49.96	ug/l	0.98	50.00	90 -	110	
65 Cu	49.87	ug/l	0.48	50.00	90 -	110	
66 Zn	50.14	ug/l	1.34	50.00	90 -	110	
75 As	48.56	ug/l	0.79	50.00	90 -	110	
78 Se	47.94	ug/l	1.13	50.00	90 -	110	
78 Se	48.27	ug/l	2.37	50.00	90 -	110	
88 Sr	50.07	ug/l	0.24	50.00	90 -	110	
88 Sr	46.85	ug/l	0.80	50.00	90 -	110	
95 Mo	48.88	ug/l	0.75	50.00	90 -	110	
106 (Cd)	--------	ug/l	--------	50.00	90 -	110	
107 Ag	24.26	ug/l	1.57	25.00	90 -	110	
108 (Cd)	--------	ug/l	--------	50.00	90 -	110	
111 Cd	49.33	ug/l	0.71	50.00	90 -	110	
118 Sn	50.24	ug/l	1.40	50.00	90 -	110	
121 Sb	49.54	ug/l	1.18	50.00	90 -	110	
137 Ba	49.63	ug/l	2.49	50.00	90 -	110	
205 Tl	48.65	ug/l	0.58	50.00	90 -	110	
206 (Pb)	--------	ug/l	--------	50.00	90 -	110	
207 (Pb)	--------	ug/l	--------	50.00	90 -	110	
208 Pb	49.96	ug/l	0.90	50.00	90 -	110	

ISTD Elements

Element	CPS Mean	RSD(%)	Ref Value	Rec(%)	QC Range(%)		Flag
6 Li	3188763.30	1.52	2775704.50	114.9	70 -	120	
45 Sc	530164.38	0.11	500780.41	105.9	70 -	120	
45 Sc	96337.35	1.21	95494.08	100.9	70 -	120	
45 Sc	1476239.00	1.18	1460980.80	101.0	70 -	120	
72 Ge	102958.30	0.45	96219.04	107.0	70 -	120	
72 Ge	45995.51	1.09	43611.78	105.5	70 -	120	
72 Ge	211979.86	0.56	213204.63	99.4	70 -	120	
115 In	1355180.90	1.54	1381264.00	98.1	70 -	120	
159 Tb	1863114.30	0.80	1843940.90	101.0	70 -	120	
165 Ho	1880561.90	0.45	1844184.90	102.0	70 -	120	

ISTD Ref File : C:\ICPCHEM\1\DATA\11K11100.B\004CALB.D\004CALB.D#

1 :Element Failures 0 :Max. Number of Failures Allowed
0 :ISTD Failures 0 :Max. Number of ISTD Failures Allowed

Data Results:
Analytes: Fail
ISTD: Pass

CCB QC Report

Data File: C:\ICPCHEM\1\DATA\11K11100.B\034_CCB.D\034_CCB.D#
Date Acquired: Nov 11 2011 03:17 pm
Operator: NBS
Sample Name: CCB 111111
Misc Info:
Vial Number: 1102
Current Method: C:\ICPCHEM\1\METHODS\62A1111A.M
Calibration File: C:\ICPCHEM\1\CALIB\62A1111A.C
Last Cal Update: Nov 11 2011 12:36 pm
Sample Type: CCB
Total Dil Factor: 1.00

QC Elements

Element	Conc.	RSD(%)	High Limit	Flag
7 (Li)	-------- ug/l	--------	#VALUE!	
9 Be	0.00 ug/l	69.30	0.12	
11 B	-0.21 ug/l	1.18	15.00	
23 Na	-12.29 ug/l	7.14	77.10	
24 Mg	0.25 ug/l	17.30	7.50	
27 Al	0.17 ug/l	52.39	3.96	
39 K	-13.38 ug/l	34.73	19.20	
44 Ca	-0.64 ug/l	345.23	90.00	
47 Ti	0.01 ug/l	249.07	0.78	
51 V	1.42 ug/l	3.15	0.21	Fail
52 Cr	0.03 ug/l	9.62	0.12	
55 Mn	0.01 ug/l	149.75	0.18	
56 Fe	0.83 ug/l	6.84	40.80	
59 Co	-0.29 ug/l	0.85	0.09	
60 Ni	-0.01 ug/l	222.72	0.48	
63 Cu	-0.38 ug/l	4.80	0.39	
65 Cu	-0.41 ug/l	3.73	0.39	
66 Zn	0.00 ug/l	2628.40	6.90	
75 As	0.07 ug/l	32.28	0.27	
78 Se	0.02 ug/l	97.30	0.30	
78 Se	0.10 ug/l	82.67	0.30	
88 Sr	0.00 ug/l	540.84	0.03	
88 Sr	0.00 ug/l	24.02	0.03	
95 Mo	0.03 ug/l	20.10	0.21	
106 (Cd)	-------- ug/l	--------	#VALUE!	
107 Ag	0.00 ug/l	72.34	0.09	
108 (Cd)	-------- ug/l	--------	#VALUE!	
111 Cd	0.00 ug/l	426.65	0.06	
118 Sn	0.05 ug/l	13.48	0.30	
121 Sb	0.29 ug/l	11.14	0.03	Fail
137 Ba	0.02 ug/l	76.66	0.12	
205 Tl	0.01 ug/l	1.73	0.03	
206 (Pb)	-------- ug/l	--------	#VALUE!	
207 (Pb)	-------- ug/l	--------	#VALUE!	
208 Pb	-0.25 ug/l	1.69	0.33	

ISTD Elements

Element	CPS Mean	RSD(%)	Ref Value	Rec(%)	QC Range(%)	Flag
6 Li	3085838.80	0.28	2775704.50	111.2	70 - 120	
45 Sc	543008.13	0.74	500780.41	108.4	70 - 120	
45 Sc	96730.10	0.69	95494.08	101.3	70 - 120	
45 Sc	1456388.00	1.14	1460980.80	99.7	70 - 120	
72 Ge	104225.84	0.21	96219.04	108.3	70 - 120	
72 Ge	45874.70	1.25	43611.78	105.2	70 - 120	
72 Ge	211968.23	0.39	213204.63	99.4	70 - 120	
115 In	1335750.50	0.67	1381264.00	96.7	70 - 120	
159 Tb	1825624.90	0.43	1843940.90	99.0	70 - 120	
165 Ho	1821355.50	0.60	1844184.90	98.8	70 - 120	

ISTD Ref File : C:\ICPCHEM\1\DATA\11K11100.B\004CALB.D\004CALB.D#

2 :Element Failures 0 :Max. Number of Failures Allowed
0 :ISTD Failures 0 :Max. Number of ISTD Failures Allowed

Data Results:
Analytes: Fail
ISTD: Pass

CCV QC Report

Data File: C:\ICPCHEM\1\DATA\11K11100.B\046_CCV.D\046_CCV.D#
Date Acquired: Nov 11 2011 04:30 pm
Operator: NBS
Sample Name: CCV 111111
Misc Info:
Vial Number: 1105
Current Method: C:\ICPCHEM\1\METHODS\62A1111A.M
Calibration File: C:\ICPCHEM\1\CALIB\62A1111A.C
Last Cal Update: Nov 11 2011 12:36 pm
Sample Type: CCV
Total Dil Factor: 1.00

QC Elements

Element	Conc.		RSD(%)	Expected	QC Range(%)		Flag
7 (Li)	--------	ug/l	--------	50.00	90 -	110	
9 Be	45.70	ug/l	1.26	50.00	90 -	110	
11 B	42.89	ug/l	1.19	50.00	90 -	110	Fail
23 Na	1254.00	ug/l	0.41	1250.00	90 -	110	
24 Mg	2547.00	ug/l	0.59	2500.00	90 -	110	
27 Al	995.00	ug/l	1.68	1000.00	90 -	110	
39 K	893.20	ug/l	1.01	1000.00	90 -	110	Fail
44 Ca	2454.00	ug/l	1.15	2500.00	90 -	110	
47 Ti	49.64	ug/l	0.77	50.00	90 -	110	
51 V	51.07	ug/l	0.56	50.00	90 -	110	
52 Cr	49.58	ug/l	1.32	50.00	90 -	110	
55 Mn	53.75	ug/l	1.23	50.00	90 -	110	
56 Fe	1008.00	ug/l	1.04	1000.00	90 -	110	
59 Co	50.27	ug/l	0.38	50.00	90 -	110	
60 Ni	50.83	ug/l	0.09	50.00	90 -	110	
63 Cu	49.08	ug/l	0.21	50.00	90 -	110	
65 Cu	49.49	ug/l	0.52	50.00	90 -	110	
66 Zn	49.79	ug/l	0.97	50.00	90 -	110	
75 As	48.81	ug/l	0.50	50.00	90 -	110	
78 Se	46.71	ug/l	1.36	50.00	90 -	110	
78 Se	47.88	ug/l	0.76	50.00	90 -	110	
88 Sr	50.09	ug/l	1.50	50.00	90 -	110	
88 Sr	46.19	ug/l	0.77	50.00	90 -	110	
95 Mo	48.00	ug/l	1.20	50.00	90 -	110	
106 (Cd)	--------	ug/l	--------	50.00	90 -	110	
107 Ag	24.50	ug/l	0.88	25.00	90 -	110	
108 (Cd)	--------	ug/l	--------	50.00	90 -	110	
111 Cd	49.59	ug/l	0.88	50.00	90 -	110	
118 Sn	50.70	ug/l	0.24	50.00	90 -	110	
121 Sb	50.92	ug/l	1.07	50.00	90 -	110	
137 Ba	50.15	ug/l	0.91	50.00	90 -	110	
205 Tl	48.73	ug/l	1.15	50.00	90 -	110	
206 (Pb)	--------	ug/l	--------	50.00	90 -	110	
207 (Pb)	--------	ug/l	--------	50.00	90 -	110	
208 Pb	50.41	ug/l	1.38	50.00	90 -	110	

ISTD Elements

Element	CPS Mean	RSD(%)	Ref Value	Rec(%)	QC Range(%)		Flag
6 Li	3087745.80	0.94	2775704.50	111.2	70 -	120	
45 Sc	518583.69	0.47	500780.41	103.6	70 -	120	
45 Sc	93007.50	0.55	95494.08	97.4	70 -	120	
45 Sc	1424009.10	1.38	1460980.80	97.5	70 -	120	
72 Ge	100723.93	0.27	96219.04	104.7	70 -	120	
72 Ge	44570.69	0.32	43611.78	102.2	70 -	120	
72 Ge	210088.66	0.52	213204.63	98.5	70 -	120	
115 In	1313527.60	0.88	1381264.00	95.1	70 -	120	
159 Tb	1801651.50	1.57	1843940.90	97.7	70 -	120	
165 Ho	1809552.10	1.53	1844184.90	98.1	70 -	120	

ISTD Ref File : C:\ICPCHEM\1\DATA\11K11100.B\004CALB.D\004CALB.D#

2 :Element Failures 0 :Max. Number of Failures Allowed
0 :ISTD Failures 0 :Max. Number of ISTD Failures Allowed

Data Results:
Analytes: Fail
ISTD: Pass

CCB QC Report

Data File: C:\ICPCHEM\1\DATA\11K11100.B\048_CCB.D\048_CCB.D#
Date Acquired: Nov 11 2011 04:42 pm
Operator: NBS
Sample Name: CCB 111111
Misc Info:
Vial Number: 1102
Current Method: C:\ICPCHEM\1\METHODS\62A1111A.M
Calibration File: C:\ICPCHEM\1\CALIB\62A1111A.C
Last Cal Update: Nov 11 2011 12:36 pm
Sample Type: CCB
Total Dil Factor: 1.00

QC Elements

Element	Conc.		RSD(%)	High Limit	Flag
7 (Li)	--------	ug/l	--------	#VALUE!	
9 Be	0.00	ug/l	42.29	0.12	
11 B	-0.13	ug/l	71.11	15.00	
23 Na	-8.31	ug/l	8.06	77.10	
24 Mg	0.36	ug/l	29.70	7.50	
27 Al	0.19	ug/l	47.04	3.96	
39 K	-21.23	ug/l	17.60	19.20	
44 Ca	-3.80	ug/l	61.16	90.00	
47 Ti	0.11	ug/l	173.80	0.78	
51 V	2.15	ug/l	1.51	0.21	Fail
52 Cr	0.08	ug/l	5.21	0.12	
55 Mn	0.54	ug/l	4.21	0.18	Fail
56 Fe	1.18	ug/l	2.51	40.80	
59 Co	-0.29	ug/l	0.62	0.09	
60 Ni	-0.01	ug/l	26.89	0.48	
63 Cu	-0.51	ug/l	3.25	0.39	
65 Cu	-0.52	ug/l	1.59	0.39	
66 Zn	0.03	ug/l	92.54	6.90	
75 As	0.35	ug/l	13.98	0.27	Fail
78 Se	0.04	ug/l	38.02	0.30	
78 Se	0.10	ug/l	95.97	0.30	
88 Sr	0.00	ug/l	152.47	0.03	
88 Sr	0.01	ug/l	57.39	0.03	
95 Mo	0.04	ug/l	33.06	0.21	
106 (Cd)	--------	ug/l	--------	#VALUE!	
107 Ag	0.00	ug/l	67.11	0.09	
108 (Cd)	--------	ug/l	--------	#VALUE!	
111 Cd	0.00	ug/l	1389.00	0.06	
118 Sn	0.07	ug/l	25.68	0.30	
121 Sb	0.41	ug/l	2.24	0.03	Fail
137 Ba	0.01	ug/l	46.25	0.12	
205 Tl	0.01	ug/l	17.61	0.03	
206 (Pb)	--------	ug/l	--------	#VALUE!	
207 (Pb)	--------	ug/l	--------	#VALUE!	
208 Pb	-0.24	ug/l	0.75	0.33	

ISTD Elements

Element	CPS Mean	RSD(%)	Ref Value	Rec(%)	QC Range(%)	Flag
6 Li	3066933.50	0.82	2775704.50	110.5	70 - 120	
45 Sc	480667.06	6.48	500780.41	96.0	70 - 120	
45 Sc	92147.45	0.95	95494.08	96.5	70 - 120	
45 Sc	1424222.60	0.94	1460980.80	97.5	70 - 120	
72 Ge	96248.01	5.46	96219.04	100.0	70 - 120	
72 Ge	44554.26	0.07	43611.78	102.2	70 - 120	
72 Ge	211003.38	0.55	213204.63	99.0	70 - 120	
115 In	1322454.80	0.90	1381264.00	95.7	70 - 120	
159 Tb	1796985.40	1.28	1843940.90	97.5	70 - 120	
165 Ho	1818211.10	0.91	1844184.90	98.6	70 - 120	

ISTD Ref File : C:\ICPCHEM\1\DATA\11K11100.B\004CALB.D\004CALB.D#

4 :Element Failures 0 :Max. Number of Failures Allowed
0 :ISTD Failures 0 :Max. Number of ISTD Failures Allowed

Data Results:
Analytes: Fail
ISTD: Pass

CCV QC Report

Data File: C:\ICPCHEM\1\DATA\11K11100.B\060_CCV.D\060_CCV.D#
Date Acquired: Nov 11 2011 06:04 pm
Operator: NBS
Sample Name: CCV 111111
Misc Info:
Vial Number: 1105
Current Method: C:\ICPCHEM\1\METHODS\62A1111A.M
Calibration File: C:\ICPCHEM\1\CALIB\62A1111A.C
Last Cal Update: Nov 11 2011 12:36 pm
Sample Type: CCV
Total Dil Factor: 1.00

QC Elements

Element	Conc.	RSD(%)	Expected	QC Range(%)	Flag
7 (Li)	-------- ug/l	--------	50.00	90 - 110	
9 Be	43.59 ug/l	0.52	50.00	90 - 110	Fail
11 B	39.89 ug/l	0.37	50.00	90 - 110	Fail
23 Na	1206.00 ug/l	0.25	1250.00	90 - 110	
24 Mg	2492.00 ug/l	1.14	2500.00	90 - 110	
27 Al	988.60 ug/l	0.65	1000.00	90 - 110	
39 K	883.40 ug/l	0.62	1000.00	90 - 110	Fail
44 Ca	2445.00 ug/l	1.17	2500.00	90 - 110	
47 Ti	49.55 ug/l	0.70	50.00	90 - 110	
51 V	52.18 ug/l	0.86	50.00	90 - 110	
52 Cr	49.10 ug/l	0.91	50.00	90 - 110	
55 Mn	53.67 ug/l	0.37	50.00	90 - 110	
56 Fe	999.40 ug/l	0.75	1000.00	90 - 110	
59 Co	49.93 ug/l	0.37	50.00	90 - 110	
60 Ni	50.18 ug/l	0.47	50.00	90 - 110	
63 Cu	48.01 ug/l	1.08	50.00	90 - 110	
65 Cu	47.95 ug/l	1.23	50.00	90 - 110	
66 Zn	48.48 ug/l	0.85	50.00	90 - 110	
75 As	48.40 ug/l	0.84	50.00	90 - 110	
78 Se	44.75 ug/l	1.93	50.00	90 - 110	Fail
78 Se	46.76 ug/l	2.40	50.00	90 - 110	
88 Sr	49.65 ug/l	0.75	50.00	90 - 110	
88 Sr	43.83 ug/l	0.41	50.00	90 - 110	Fail
95 Mo	45.84 ug/l	0.72	50.00	90 - 110	
106 (Cd)	-------- ug/l	--------	50.00	90 - 110	
107 Ag	24.47 ug/l	1.71	25.00	90 - 110	
108 (Cd)	-------- ug/l	--------	50.00	90 - 110	
111 Cd	48.94 ug/l	2.42	50.00	90 - 110	
118 Sn	49.84 ug/l	1.46	50.00	90 - 110	
121 Sb	50.04 ug/l	0.62	50.00	90 - 110	
137 Ba	50.55 ug/l	2.58	50.00	90 - 110	
205 Tl	46.49 ug/l	1.08	50.00	90 - 110	
206 (Pb)	-------- ug/l	--------	50.00	90 - 110	
207 (Pb)	-------- ug/l	--------	50.00	90 - 110	
208 Pb	48.04 ug/l	0.60	50.00	90 - 110	

ISTD Elements

Element	CPS Mean	RSD(%)	Ref Value	Rec(%)	QC Range(%)	Flag
6 Li	3120885.30	0.68	2775704.50	112.4	70 - 120	
45 Sc	528291.50	1.60	500780.41	105.5	70 - 120	
45 Sc	94943.62	0.72	95494.08	99.4	70 - 120	
45 Sc	1497531.60	0.60	1460980.80	102.5	70 - 120	
72 Ge	107482.91	1.52	96219.04	111.7	70 - 120	
72 Ge	46381.07	0.48	43611.78	106.3	70 - 120	
72 Ge	233866.19	0.20	213204.63	109.7	70 - 120	
115 In	1403864.60	1.34	1381264.00	101.6	70 - 120	
159 Tb	1927869.00	0.42	1843940.90	104.6	70 - 120	
165 Ho	1902582.90	0.32	1844184.90	103.2	70 - 120	

ISTD Ref File : C:\ICPCHEM\1\DATA\11K11100.B\004CALB.D\004CALB.D#

5 :Element Failures 0 :Max. Number of Failures Allowed
0 :ISTD Failures 0 :Max. Number of ISTD Failures Allowed

Data Results:
Analytes: Fail
ISTD: Pass

CCB QC Report

Data File: C:\ICPCHEM\1\DATA\11K11100.B\062_CCB.D\062_CCB.D#
Date Acquired: Nov 11 2011 06:16 pm
Operator: NBS
Sample Name: CCB 111111
Misc Info:
Vial Number: 1102
Current Method: C:\ICPCHEM\1\METHODS\62A1111A.M
Calibration File: C:\ICPCHEM\1\CALIB\62A1111A.C
Last Cal Update: Nov 11 2011 12:36 pm
Sample Type: CCB
Total Dil Factor: 1.00

QC Elements

Element	Conc.	RSD(%)	High Limit	Flag
7 (Li)	-------- ug/l	--------	#VALUE!	
9 Be	0.00 ug/l	45.82	0.12	
11 B	-0.52 ug/l	3.32	15.00	
23 Na	-14.13 ug/l	1.97	77.10	
24 Mg	0.40 ug/l	18.80	7.50	
27 Al	0.37 ug/l	36.41	3.96	
39 K	-25.01 ug/l	22.10	19.20	
44 Ca	-5.06 ug/l	39.41	90.00	
47 Ti	-0.01 ug/l	153.32	0.78	
51 V	3.23 ug/l	1.14	0.21	Fail
52 Cr	0.11 ug/l	18.29	0.12	
55 Mn	0.56 ug/l	2.13	0.18	Fail
56 Fe	0.91 ug/l	1.90	40.80	
59 Co	-0.29 ug/l	0.95	0.09	
60 Ni	-0.01 ug/l	26.61	0.48	
63 Cu	-0.57 ug/l	1.33	0.39	
65 Cu	-0.57 ug/l	1.17	0.39	
66 Zn	0.10 ug/l	36.98	6.90	
75 As	0.63 ug/l	2.54	0.27	Fail
78 Se	0.03 ug/l	36.72	0.30	
78 Se	0.20 ug/l	17.06	0.30	
88 Sr	0.01 ug/l	140.22	0.03	
88 Sr	0.00 ug/l	12.28	0.03	
95 Mo	0.01 ug/l	62.46	0.21	
106 (Cd)	-------- ug/l	--------	#VALUE!	
107 Ag	0.00 ug/l	55.01	0.09	
108 (Cd)	-------- ug/l	--------	#VALUE!	
111 Cd	0.01 ug/l	225.87	0.06	
118 Sn	0.05 ug/l	34.43	0.30	
121 Sb	0.19 ug/l	1.47	0.03	Fail
137 Ba	0.00 ug/l	114.50	0.12	
205 Tl	0.01 ug/l	8.77	0.03	
206 (Pb)	-------- ug/l	--------	#VALUE!	
207 (Pb)	-------- ug/l	--------	#VALUE!	
208 Pb	-0.25 ug/l	1.47	0.33	

ISTD Elements

Element	CPS Mean	RSD(%)	Ref Value	Rec(%)	QC Range(%)	Flag
6 Li	3025901.00	0.96	2775704.50	109.0	70 - 120	
45 Sc	540897.69	0.43	500780.41	108.0	70 - 120	
45 Sc	95060.94	0.12	95494.08	99.5	70 - 120	
45 Sc	1475771.40	0.55	1460980.80	101.0	70 - 120	
72 Ge	108235.30	0.84	96219.04	112.5	70 - 120	
72 Ge	46007.31	1.05	43611.78	105.5	70 - 120	
72 Ge	232509.75	0.78	213204.63	109.1	70 - 120	
115 In	1409864.80	1.11	1381264.00	102.1	70 - 120	
159 Tb	1904300.90	0.32	1843940.90	103.3	70 - 120	
165 Ho	1879356.80	0.14	1844184.90	101.9	70 - 120	

ISTD Ref File : C:\ICPCHEM\1\DATA\11K11100.B\004CALB.D\004CALB.D#

4 :Element Failures 0 :Max. Number of Failures Allowed
0 :ISTD Failures 0 :Max. Number of ISTD Failures Allowed

Data Results:
Analytes: Fail
ISTD: Pass

METALS
Raw Data

APPL, INC.

METALS BLANK

APPL Inc.
908 North Temperance Avenue
Clovis, CA 93611

Method	Analyte	Result	LOQ	LOD	DL	Units	Prep Date	Analysis Date	QC Group
6020	LEAD (PB) (DISSOL	0.19 J	0.5	0.22	0.11	ug/L	11/10/11	11/11/11	#602D-111110A-AY49334

J = Estimated value.

Metals SC-Blank-REG MDLs
Printed: 11/15/2011 12:58:31 PM

Sample QC Report

Data File: C:\ICPCHEM\1\DATA\11K11100.B\024SMPL.D\024SMPL.D#
Date Acquired: Nov 11 2011 02:16 pm
Operator: NBS
Sample Name: 111110A-3015-BLK
Misc Info: 111110A-3015
Vial Number: 3101
Current Method: C:\ICPCHEM\1\METHODS\62A1111A.M
Calibration File: C:\ICPCHEM\1\CALIB\62A1111A.C
Last Cal Update: Nov 11 2011 12:36 pm
Sample Type: Sample
Prep Dil Factor: 1.11
Total Dil Factor: 1.11

QC Elements

Element	Conc.		Corr. Conc.	RSD(%)	High Limit	Flag
7 (Li)	--------	ug/l	#VALUE!	--------	0	
9 Be	-0.01	ug/l	-0.01	2.73	1000	
11 B	0.02	ug/l	0.02	15.56	1000	
23 Na	35.01	ug/l	38.90	12.71	25000	
24 Mg	5.25	ug/l	5.83	2.15	50000	
27 Al	6.63	ug/l	7.37	3.21	20000	
39 K	-19.19	ug/l	-21.32	24.75	20000	
44 Ca	187.10	ug/l	207.87	2.16	50000	
47 Ti	0.09	ug/l	0.10	43.80	1000	
51 V	-0.78	ug/l	-0.86	1.87	1000	
52 Cr	-0.04	ug/l	-0.04	28.95	1000	
55 Mn	0.23	ug/l	0.26	6.88	1000	
56 Fe	2.70	ug/l	3.00	5.14	20000	
59 Co	-0.27	ug/l	-0.30	0.86	1000	
60 Ni	0.12	ug/l	0.14	23.87	1000	
63 Cu	-0.44	ug/l	-0.49	2.50	1000	
65 Cu	-0.44	ug/l	-0.49	2.58	1000	
66 Zn	7.48	ug/l	8.31	3.75	1000	
75 As	-0.53	ug/l	-0.59	2.55	1000	
78 Se	-0.01	ug/l	-0.01	26.71	1000	
78 Se	-0.01	ug/l	-0.02	520.99	1000	
88 Sr	0.14	ug/l	0.16	12.74	1000	
88 Sr	0.14	ug/l	0.16	3.88	1000	
95 Mo	0.02	ug/l	0.02	6.80	1000	
106 (Cd)	--------	ug/l	#VALUE!	--------	########	
107 Ag	0.00	ug/l	0.00	212.49	500	
108 (Cd)	--------	ug/l	#VALUE!	--------	########	
111 Cd	0.02	ug/l	0.02	26.66	1000	
118 Sn	0.12	ug/l	0.13	6.01	1000	
121 Sb	0.07	ug/l	0.08	7.73	1000	
137 Ba	0.04	ug/l	0.04	9.41	1000	
205 Tl	0.01	ug/l	0.01	22.33	1000	
206 (Pb)	--------	ug/l	#VALUE!	--------	########	
207 (Pb)	--------	ug/l	#VALUE!	--------	########	
208 Pb	0.17	ug/l	0.19	1.81	1000	

ISTD Elements

Element	CPS Mean	RSD(%)	Ref Value	Rec(%)	QC Range(%)	Flag
6 Li	3205542.30	0.83	2775704.50	115.5	70 - 120	
45 Sc	579022.81	0.89	500780.41	115.6	70 - 120	
45 Sc	106222.45	0.45	95494.08	111.2	70 - 120	
45 Sc	1635333.40	0.66	1460980.80	111.9	70 - 120	
72 Ge	108091.34	0.67	96219.04	112.3	70 - 120	
72 Ge	49642.59	1.23	43611.78	113.8	70 - 120	
72 Ge	228973.69	0.34	213204.63	107.4	70 - 120	
115 In	1505106.90	0.63	1381264.00	109.0	70 - 120	
159 Tb	2069441.80	1.02	1843940.90	112.2	70 - 120	
165 Ho	2056674.30	0.60	1844184.90	111.5	70 - 120	

ISTD Ref File : C:\ICPCHEM\1\DATA\11K11100.B\004CALB.D\004CALB.D#

0 :Element Failures 0 :Max. Number of Failures Allowed
0 :ISTD Failures 0 :Max. Number of ISTD Failures Allowed

Data Results:
Analytes: Pass
ISTD: Pass

Laboratory Control Spike Recovery

METALS

APPL Inc.
908 North Temperance Avenue
Clovis, CA 93611

Method	Compound Name	Spike Level ug/L	SPK Result ug/L	SPK % Recovery	Recovery Limits	Extract Date	Analysis Date	QC Group
6020	LEAD (PB) (DISSOLVED)	50.0	50.0	100	80-120	11/10/2011	1/11/2011	#602D-111110A-AY49334

Comments: _____

Sample QC Report

Data File: C:\ICPCHEM\1\DATA\11K11100.B\025SMPL.D\025SMPL.D#
Date Acquired: Nov 11 2011 02:22 pm
Operator: NBS
Sample Name: 111110A-3015-LCS
Misc Info: 111110A-3015
Vial Number: 3102
Current Method: C:\ICPCHEM\1\METHODS\62A1111A.M
Calibration File: C:\ICPCHEM\1\CALIB\62A1111A.C
Last Cal Update: Nov 11 2011 12:36 pm
Sample Type: Sample
Prep Dil Factor: 1.11
Total Dil Factor: 1.11

QC Elements

Element	Conc.		Corr. Conc.	RSD(%)	High Limit	Flag
7 (Li)	--------	ug/l	#VALUE!	--------	0	
9 Be	7.09	ug/l	7.88	0.95	1000	
11 B	34.46	ug/l	38.29	0.83	1000	
23 Na	4313.00	ug/l	4791.74	0.33	25000	
24 Mg	4287.00	ug/l	4762.86	0.50	50000	
27 Al	378.80	ug/l	420.85	1.07	20000	
39 K	819.80	ug/l	910.80	1.14	20000	
44 Ca	4772.00	ug/l	5301.69	0.81	50000	
47 Ti	43.16	ug/l	47.95	0.73	1000	
51 V	44.89	ug/l	49.87	0.68	1000	
52 Cr	46.85	ug/l	52.05	0.52	1000	
55 Mn	48.25	ug/l	53.61	0.34	1000	
56 Fe	188.60	ug/l	209.53	0.62	20000	
59 Co	45.14	ug/l	50.15	0.81	1000	
60 Ni	45.24	ug/l	50.26	0.75	1000	
63 Cu	42.57	ug/l	47.30	0.51	1000	
65 Cu	42.70	ug/l	47.44	0.10	1000	
66 Zn	94.53	ug/l	105.02	0.68	1000	
75 As	39.81	ug/l	44.23	0.66	1000	
78 Se	36.57	ug/l	40.63	2.90	1000	
78 Se	37.79	ug/l	41.98	1.41	1000	
88 Sr	47.19	ug/l	52.43	0.39	1000	
88 Sr	45.26	ug/l	50.28	0.18	1000	
95 Mo	45.43	ug/l	50.47	0.63	1000	
106 (Cd)	--------	ug/l	#VALUE!	--------	########	
107 Ag	16.57	ug/l	18.41	1.13	500	
108 (Cd)	--------	ug/l	#VALUE!	--------	########	
111 Cd	8.34	ug/l	9.26	2.50	1000	
118 Sn	48.01	ug/l	53.34	0.41	1000	
121 Sb	42.84	ug/l	47.60	0.50	1000	
137 Ba	44.60	ug/l	49.55	0.91	1000	
205 Tl	43.45	ug/l	48.27	0.08	1000	
206 (Pb)	--------	ug/l	#VALUE!	--------	########	
207 (Pb)	--------	ug/l	#VALUE!	--------	########	
208 Pb	45.08	ug/l	50.08	0.48	1000	

ISTD Elements

Element	CPS Mean	RSD(%)	Ref Value	Rec(%)	QC Range(%)	Flag
6 Li	3106113.00	0.94	2775704.50	111.9	70 - 120	
45 Sc	583837.94	0.76	500780.41	116.6	70 - 120	
45 Sc	104815.63	0.97	95494.08	109.8	70 - 120	
45 Sc	1623628.90	0.72	1460980.80	111.1	70 - 120	
72 Ge	109519.99	0.75	96219.04	113.8	70 - 120	
72 Ge	48705.67	0.97	43611.78	111.7	70 - 120	
72 Ge	226177.02	0.48	213204.63	106.1	70 - 120	
115 In	1499201.30	0.50	1381264.00	108.5	70 - 120	
159 Tb	2052386.10	0.41	1843940.90	111.3	70 - 120	
165 Ho	2061841.80	0.62	1844184.90	111.8	70 - 120	

ISTD Ref File : C:\ICPCHEM\1\DATA\11K11100.B\004CALB.D\004CALB.D#

0 :Element Failures 0 :Max. Number of Failures Allowed
0 :ISTD Failures 0 :Max. Number of ISTD Failures Allowed

Data Results:
Analytes: Pass
ISTD: Pass

Matrix Spike Recoveries

METALS

APPL ID: **111110W-49334 MS - 161255**

Sample ID: AY49334

Client ID: ES047

APPL Inc.
908 North Temperance Avenue
Clovis, CA 93611

Method	Compound Name	Spike Lvl ug/L	Matrix Res ug/L	SPK Res ug/L	DUP Res ug/L	SPK % Recovery	DUP % Recovery	RPD	RPD Max	Recovery Limits	Extract Date-Spk	Analysis Date-Spk	Extract Date-Dup	Analysis Date-Dup	QC Group	QC Sample
6020	LEAD (PB) (DISSOLVE	50.0	ND	49.5	48.8	99.0	97.6	1.4	20	80-120	1/10/2011	1/11/2011	1/10/2011	1/11/2011	161255	AY49334

Comments: _____

Sample QC Report

Data File: C:\ICPCHEM\1\DATA\11K11100.B\043SMPL.D\043SMPL.D#
Date Acquired: Nov 11 2011 04:12 pm
Operator: NBS
Sample Name: AY49334W52 MS
Misc Info: 111110A-3015
Vial Number: 3204
Current Method: C:\ICPCHEM\1\METHODS\62A1111A.M
Calibration File: C:\ICPCHEM\1\CALIB\62A1111A.C
Last Cal Update: Nov 11 2011 12:36 pm
Sample Type: Sample
Prep Dil Factor: 1.11
Total Dil Factor: 1.11

QC Elements

Element	Conc.		Corr. Conc.	RSD(%)	High Limit	Flag
7 (Li)	--------	ug/l	#VALUE!	--------	0	
9 Be	5.49	ug/l	6.10	1.06	1000	
11 B	106.30	ug/l	118.10	0.86	1000	
23 Na	67250.00	ug/l	74714.75	0.69	25000	>Cal
24 Mg	27980.00	ug/l	31085.78	0.94	50000	
27 Al	425.20	ug/l	472.40	1.20	20000	
39 K	3527.00	ug/l	3918.50	0.62	20000	
44 Ca	17630.00	ug/l	19586.93	1.01	50000	
47 Ti	46.43	ug/l	51.58	0.68	1000	
51 V	47.50	ug/l	52.77	0.75	1000	
52 Cr	46.44	ug/l	51.59	0.93	1000	
55 Mn	1471.00	ug/l	1634.28	0.54	1000	>Cal
56 Fe	1106.00	ug/l	1228.77	0.88	20000	
59 Co	44.68	ug/l	49.64	1.20	1000	
60 Ni	44.83	ug/l	49.81	0.65	1000	
63 Cu	43.23	ug/l	48.03	0.77	1000	
65 Cu	43.18	ug/l	47.97	0.92	1000	
66 Zn	91.21	ug/l	101.33	0.62	1000	
75 As	42.58	ug/l	47.31	0.67	1000	
78 Se	38.49	ug/l	42.76	2.21	1000	
78 Se	39.09	ug/l	43.43	0.75	1000	
88 Sr	163.00	ug/l	181.09	0.38	1000	
88 Sr	163.50	ug/l	181.65	1.22	1000	
95 Mo	47.02	ug/l	52.24	1.05	1000	
106 (Cd)	--------	ug/l	#VALUE!	--------	########	
107 Ag	17.43	ug/l	19.36	0.62	500	
108 (Cd)	--------	ug/l	#VALUE!	--------	########	
111 Cd	8.85	ug/l	9.83	1.40	1000	
118 Sn	49.73	ug/l	55.25	0.73	1000	
121 Sb	43.10	ug/l	47.88	0.46	1000	
137 Ba	71.28	ug/l	79.19	1.16	1000	
205 Tl	43.09	ug/l	47.87	0.35	1000	
206 (Pb)	--------	ug/l	#VALUE!	--------	########	
207 (Pb)	--------	ug/l	#VALUE!	--------	########	
208 Pb	44.63	ug/l	49.58	1.25	1000	

ISTD Elements

Element	CPS Mean	RSD(%)	Ref Value	Rec(%)	QC Range(%)	Flag
6 Li	3108599.50	0.69	2775704.50	112.0	70 - 120	
45 Sc	657553.31	0.13	500780.41	131.3	70 - 120	IS Fai
45 Sc	107348.46	1.08	95494.08	112.4	70 - 120	
45 Sc	2219467.30	0.83	1460980.80	151.9	70 - 120	IS Fai
72 Ge	104662.85	0.47	96219.04	108.8	70 - 120	
72 Ge	48176.50	1.42	43611.78	110.5	70 - 120	
72 Ge	232443.67	0.24	213204.63	109.0	70 - 120	
115 In	1472805.40	0.82	1381264.00	106.6	70 - 120	
159 Tb	2035764.30	0.83	1843940.90	110.4	70 - 120	
165 Ho	2055214.00	0.90	1844184.90	111.4	70 - 120	

ISTD Ref File : C:\ICPCHEM\1\DATA\11K11100.B\004CALB.D\004CALB.D#

2 :Element Failures 0 :Max. Number of Failures Allowed
2 :ISTD Failures 0 :Max. Number of ISTD Failures Allowed

Data Results:
Analytes: Fail
ISTD: Fail

Sample QC Report

Data File: C:\ICPCHEM\1\DATA\11K11100.B\049SMPL.D\049SMPL.D#
Date Acquired: Nov 11 2011 04:48 pm
Operator: NBS
Sample Name: AY49334WS2 MSD
Misc Info: 111110A-3015
Vial Number: 3205
Current Method: C:\ICPCHEM\1\METHODS\62A1111A.M
Calibration File: C:\ICPCHEM\1\CALIB\62A1111A.C
Last Cal Update: Nov 11 2011 12:36 pm
Sample Type: Sample
Prep Dil Factor: 1.11
Total Dil Factor: 1.11

QC Elements

Element	Conc.		Corr. Conc.	RSD(%)	High Limit	Flag
7 (Li)	--------	ug/l	#VALUE!	--------	0	
9 Be	5.56	ug/l	6.17	0.61	1000	
11 B	105.40	ug/l	117.10	0.69	1000	
23 Na	67550.00	ug/l	75048.05	0.54	25000	>Cal
24 Mg	27890.00	ug/l	30985.79	0.29	50000	
27 Al	375.40	ug/l	417.07	1.44	20000	
39 K	3571.00	ug/l	3967.38	0.91	20000	
44 Ca	17520.00	ug/l	19464.72	1.41	50000	
47 Ti	46.01	ug/l	51.12	3.10	1000	
51 V	48.57	ug/l	53.96	1.00	1000	
52 Cr	47.16	ug/l	52.39	1.21	1000	
55 Mn	1476.00	ug/l	1639.84	0.85	1000	>Cal
56 Fe	1047.00	ug/l	1163.22	0.58	20000	
59 Co	45.69	ug/l	50.76	0.38	1000	
60 Ni	45.51	ug/l	50.56	0.70	1000	
63 Cu	43.37	ug/l	48.18	0.83	1000	
65 Cu	43.50	ug/l	48.33	1.42	1000	
66 Zn	81.55	ug/l	90.60	0.72	1000	
75 As	43.24	ug/l	48.04	0.69	1000	
78 Se	39.39	ug/l	43.76	2.26	1000	
78 Se	40.07	ug/l	44.52	0.71	1000	
88 Sr	163.50	ug/l	181.65	0.57	1000	
88 Sr	165.60	ug/l	183.98	0.32	1000	
95 Mo	48.36	ug/l	53.73	0.93	1000	
106 (Cd)	--------	ug/l	#VALUE!	--------	#######	
107 Ag	17.97	ug/l	19.96	1.57	500	
108 (Cd)	--------	ug/l	#VALUE!	--------	#######	
111 Cd	8.84	ug/l	9.82	0.94	1000	
118 Sn	50.45	ug/l	56.05	1.04	1000	
121 Sb	46.12	ug/l	51.24	0.92	1000	
137 Ba	70.84	ug/l	78.70	1.10	1000	
205 Tl	42.97	ug/l	47.74	0.67	1000	
206 (Pb)	--------	ug/l	#VALUE!	--------	#######	
207 (Pb)	--------	ug/l	#VALUE!	--------	#######	
208 Pb	43.94	ug/l	48.82	1.12	1000	

ISTD Elements

Element	CPS Mean	RSD(%)	Ref Value	Rec(%)	QC Range(%)	Flag
6 Li	3085573.50	1.49	2775704.50	111.2	70 - 120	
45 Sc	659528.88	0.85	500780.41	131.7	70 - 120	IS Fai
45 Sc	103045.15	1.13	95494.08	107.9	70 - 120	
45 Sc	2177886.80	0.53	1460980.80	149.1	70 - 120	IS Fai
72 Ge	102879.23	0.86	96219.04	106.9	70 - 120	
72 Ge	46907.60	1.75	43611.78	107.6	70 - 120	
72 Ge	224264.44	0.97	213204.63	105.2	70 - 120	
115 In	1453996.00	1.44	1381264.00	105.3	70 - 120	
159 Tb	2021027.00	0.98	1843940.90	109.6	70 - 120	
165 Ho	2010739.50	0.60	1844184.90	109.0	70 - 120	

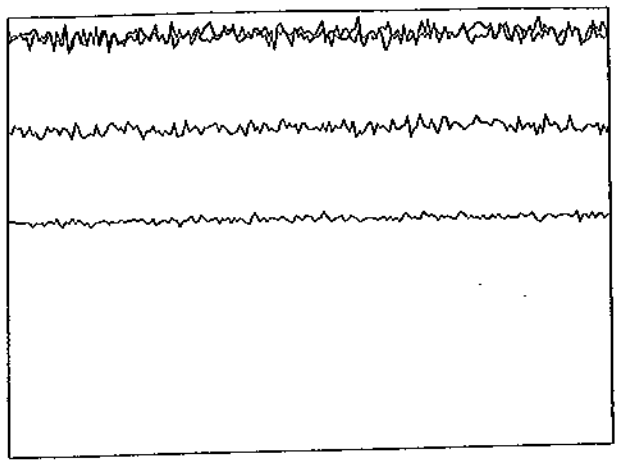
ISTD Ref File : C:\ICPCHEM\1\DATA\11K11100.B\004CALB.D\004CALB.D#

2 :Element Failures 0 :Max. Number of Failures Allowed
2 :ISTD Failures 0 :Max. Number of ISTD Failures Allowed

Data Results:
Analytes: Fail
ISTD: Fail

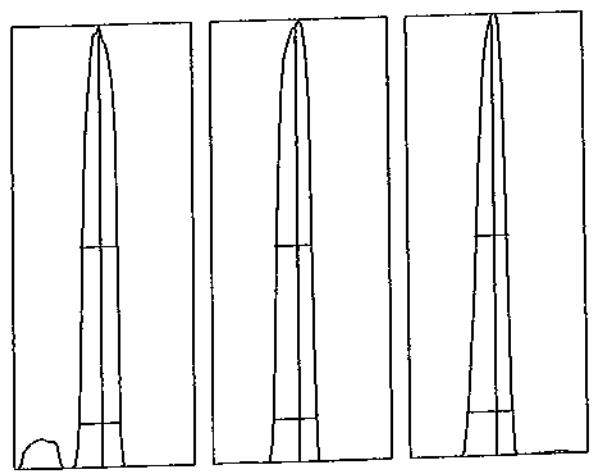
Tune Report

Tune File : nogas.u
Comment : 111111



Integration Time:	0.1000 sec	
Sampling Period:	0.6200 sec	
n:	200	
Oxide:	156/140	1.410%
Doubly Charged:	70/140	1.051%

m/z	Range	Count	Mean	RSD%	Background
7	50,000	26283.0	26440.3	1.09	0.40
89	20,000	19274.0	18861.9	1.39	2.20
205	20,000	14914.0	14722.1	1.50	5.80
156/140	2	1.520%	1.398%	6.48	
70/140	2	1.065%	1.038%	8.17	
140	20,000	18882.0	19064.3	1.33	4.10



m/z:	7	89	205
Height:	26,611	18,699	14,936
Axis:	7.00	89.00	205.00
W-50%:	0.65	0.65	0.60
W-10%:	0.700	0.7500	0.800

Integration Time: 0.1000 sec
Acquisition Time: 22.7600 sec

Y axis : Linear

Tune Report

Tune File : nogas.u
Comment : 111111

Tuning Parameters

===Plasma Condition===			===Ion Lenses===			===Q-Pole Parameters===		
RF Power	:	1600 W	Extract 1	:	0 V	AMU Gain	:	128
RF Matching	:	1.66 V	Extract 2	:	-130 V	AMU Offset	:	127
Smpl Depth	:	9.6 mm	Omega Bias-ce	:	-22 V	Axis Gain	:	1
Torch-H	:	-0.1 mm	Omega Lens-ce	:	-1.2 V	Axis Offset	:	-0.02
Torch-V	:	0.1 mm	Cell Entrance	:	-30 V	QP Bias	:	-3 V
Carrier Gas	:	1.02 L/min	QP Focus	:	5 V			
Makeup Gas	:	0.1 L/min	Cell Exit	:	-30 V	===Detector Parameters===		
Optional Gas	:	--- %				Discriminator	:	8 mV
Nebulizer Pump	:	0.1 rps	===Octopole Parameters===			Analog HV	:	1660 V
Sample Pump	:	--- rps	OctP RF	:	180 V	Pulse HV	:	1460 V
S/C Temp	:	2 degC	OctP Bias	:	-6 V			

===Reaction Cell===

Reaction Mode	:	OFF						
H2 Gas	:	0 mL/min	He Gas	:	0 mL/min	Optional Gas	:	--- %

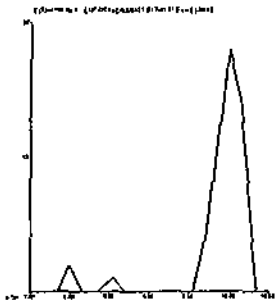
Generated : Nov 11, 2011 10:05:38
Printed : Nov 11, 2011 10:05:42

200.8 QC Tune Report

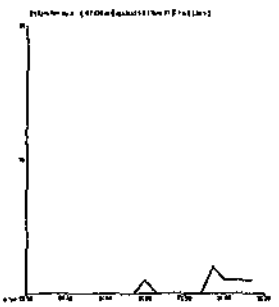
Data File: C:\ICPCHEM\1\DATA\11K11100.B\001TUNE.D
Date Acquired: Nov 11 2011 11:48 am
Acq. Method: TN200_8.M
Operator: NBS
Sample Name: 100ppb Tune sol
Misc Info:
Vial Number: 1303
Current Method: C:\ICPCHEM\1\METHODS\TN200_8.M

RSD (%)

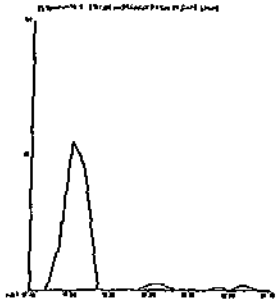
Element	CPS Mean	Rep1	Rep2	Rep3	Rep4	Rep5	%RSD	Required	Flag
9 Be	65891175	64840372	65630536	65890148	66486284	66608536	1.01	5.00	
24 Mg	120432836	########	########	########	########	########	1.16	5.00	
59 Co	111175066	########	########	########	########	########	0.73	5.00	
115 In	122240964	########	########	########	########	########	0.81	5.00	
208 Pb	63959189	64419004	64182972	63372424	64206080	63615464	1.13	5.00	



9 Be
Mass Calib.
Actual: 9.00
Required: 8.90 - 9.10
Flag:
Peak Width
Actual: 0.60
Required: 0.90
Flag:



24 Mg
Mass Calib.
Actual: 23.95
Required: 23.90 - 24.10
Flag:
Peak Width
Actual: 0.65
Required: 0.80
Flag:



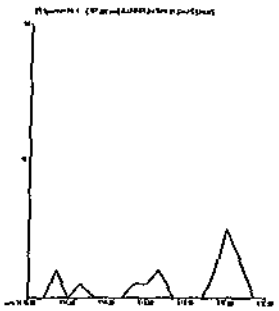
59 Co

Mass Calib.

Actual:	59.00		
Required:	58.90	-	59.10
Flag:			

Peak Width

Actual:	0.60		
Required:	0.90		
Flag:			



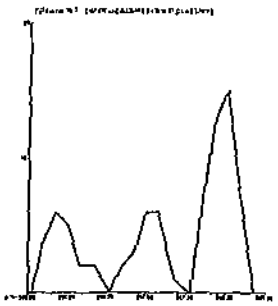
115 In

Mass Calib.

Actual:	115.05		
Required:	114.90	-	115.10
Flag:			

Peak Width

Actual:	0.65		
Required:	0.90		
Flag:			



208 Pb

Mass Calib.

Actual:	208.00		
Required:	207.90	-	208.10
Flag:			

Peak Width

Actual:	0.65		
Required:	0.80		
Flag:			

Tune Result: Pass

Metals Standards Log Book # 34 Page #001

NBS 11/11/11

NBS 11/11/11
6020/6020A
(A)

ICP-MS STANDARDS 6020/6020A/3015/3051A
Today's Date: 11/11/2011
Expires: 11/18/2011
Prep Date 1% HNO3/1.0%HCL
20 mL HNO3 / 2000 mL DI Water
Lot # K19023
20mL HCL / 2000mL DI Water
Lot #4110110
Expires: 11/18/2011

Standard 6

Amount	STD	Manufacturer	Lot #
50 uL	CCV-A	Env. Express	1036407-28139
50 uL	CCV-B	Env. Express	1036410-28140
50 uL	CCV-C	Env. Express	1100309-28141

Prepared in 100 mL of 1% HNO3/1.0% HCL 11/11/2011

Standard 3 11/18/2011

Amount	STD	Manufacturer	Lot #
25 uL	CCV-A	Env. Express	1036407-28139
25 uL	CCV-B	Env. Express	1036410-28140
25 uL	CCV-C	Env. Express	1100309-28141

Prepared in 100 mL of 1% HNO3/1.0% HCL 11/11/2011

Standard 2 11/18/2011
Amount STD
500 uL Standard 4 11/11/2011
Prepared in 50 mL of 1% HNO3/1.0% HCL 11/11/2011

Standard 1 11/18/2011
Amount STD
50 uL Standard 4 11/11/2011
Prepared in 50 mL of 1% HNO3/1.0% HCL 11/11/2011

ICP-MS ICV 11/18/2011
Amount STD

50 uL	QCS ICV A	CPI	11C174-28548
50 uL	QCS ICV B	CPI	11C174-28549

Prepared in 50 mL of 1% HNO3/1.0% HCL 11/11/2011

ICSA Prep: 11/18/2011

1 mL	ICSA	CPI	11C066-28528

Prepared in 5 mL of 1% HNO3/1.0% HCL 11/11/2011

ICSAB Prep: 11/18/2011

1mL	ICSA	CPI	11C066-28528
0.025mL	INT	O2SI	1023805-28210

Prepared in 5 mL of 1% HNO3/1.0% HCL 11/11/2011

ICP-LDR 11/18/2011
Amount STD

50 uL	CCV-A	Env. Express	1036407-28139
50 uL	CCV-B	Env. Express	1036410-28140
50 uL	CCV-C	Env. Express	1100309-28141

Prepared in 10 mL of 1% HNO3/1.0% HCL 11/11/2011

SDM 11/11/11

SDM 11/11/11
200.7
Exp (A)

2%HNO3 / 2%HCl BLK				
AMOUNT	REAGENT	MANUFACTURER	LOT	OPEN DATE
40 mL	HCL	BDH	4110110	10/14/2011
40 mL	HNO3	JT BAKER	K19023	10/14/2011
Prepared in 2000 ml DI Water				

STD 1 / LDL 200.7				
AMOUNT	STD	MANUFACTURER	LOT	EXP DATE
0.250 mL	200.7 LDL	O2SI	1028857-29687	11/1/2012
Prepared in 50 ml 2%HNO3/2%HCl				

STD 3 / HDL 200.7				
0.5 mL	CCV-A	ABSOLUTE	091409-25206	9/14/2012
0.5 mL	CCV-B	ABSOLUTE	091109-25206	9/14/2012
0.5 mL	CCV-C	ABSOLUTE	091009-25207	9/10/2012

STD 2 / CCV1 200.7				
AMOUNT	STD	PREP DATE	EXP DATE	
25mL	STD 3	11/4/2011	11/11/2011	
25mL	2%HNO3/2%HCl	11/4/2011	11/11/2011	

CCV2 200.7			
15mL	STD 3	11/4/2011	11/11/2011
25mL	2%HNO3/2%HCl	11/4/2011	11/11/2011

200.7 ICV				
AMOUNT	STD	MANUFACTURER	LOT	EXP DATE
0.5ML	QCS ICV A	CPI	11C174-28548	9/17/2012
0.5ML	QCS ICV B	CPI	11C174-28549	9/17/2012
Prepared in 50ml 2%HNO3/2%HCl				

200.7 ICSA				
0.5mL	Al	CPI	10E012-27685	4/20/2012
0.5mL	Ca	CPI	11A008-28528	9/15/2012
0.5mL	Mg	CPI	10H213-2786	4/20/2012
0.5mL	Fe	O2SI	1022245-27699	4/22/2012
Prepared in 50 ml 2%HNO3/2%HCl				

200.7 ICSAB				
0.5mL	Al	CPI	10E012-27685	4/20/2012
0.5mL	Ca	CPI	11A008-28528	9/15/2012
0.5mL	Mg	CPI	10H213-2786	4/20/2012
0.5mL	Fe	O2SI	1022245-27699	4/22/2012
0.25mL	INT SPECIAL MIX	O2SI	160495-01-01	3/1/2012
Prepared in 50 ml 2%HNO3/2%HCl				

SDM 11/11/11
6010B/6010C
(A)

1%HNO3 / 5%HCl BLK				
AMOUNT	REAGENT	MANUFACTURER	LOT	OPEN DATE
100 mL	HCL	BDH	4110110	10/14/2011
20 mL	HNO3	JT BAKER	K19023	10/14/2011
Prepared in 2000 ml DI Water				

STD 1 / LDL 6010B/6010C				
AMOUNT	STD	MANUFACTURER	LOT	EXP DATE
0.5 mL	6010 LDL	ABSOLUTE	091409-25205	9/14/2012
Prepared in 50 ml 1%HNO3/5%HCl				

STD 3 / HDL 6010B/6010C				
1ML	CCV-A	ABSOLUTE	091409-25206	9/14/2012
1ML	CCV-B	ABSOLUTE	091109-25206	9/14/2012
1ML	CCV-C	ABSOLUTE	091009-25207	9/10/2012
Prepared in 100 ml 1%HNO3 / 5%HCl				

STD 2 / CCV1 6010B/6010C/6010C				
AMOUNT	STD	PREP DATE	EXP DATE	
25mL	STD 3	11/11/2011	11/18/2011	
25mL	1%HNO3/5%HCl	11/11/2011	11/18/2011	

CCV2 6010B/6010C			
AMOUNT	STD	PREP DATE	EXP DATE
15mL	STD 3	11/11/2011	11/18/2011
25mL	1%HNO3/5%HCl	11/11/2011	11/18/2011

6010B/6010C ICSA				
AMOUNT	STD	MANUFACTURER	LOT	EXP DATE
1mL	Al	CPI	10E012-27685	4/20/2012
1mL	Ca	CPI	11A008-28528	9/15/2012
1mL	Mg	CPI	10H213-2786	4/20/2012
1mL	Fe	O2SI	1022245-27699	4/22/2012
Prepared in 50 ml 1%HNO3/5%HCl				

6010B/6010C ICSAB				
1mL	Al	CPI	10E012-27685	4/20/2012
1mL	Ca	CPI	11A008-28528	9/15/2012
1mL	Mg	CPI	10H213-2786	4/20/2012
1mL	Fe	O2SI	1022245-27699	4/22/2012
0.5mL	JT SPECIAL M	O2SI	160495-01-01	3/1/2012
Prepared in 50 ml 1%HNO3/5%HCl				

6010B/6010C ICV				
0.5ML	QCS ICV A	CPI	11C174-28548	9/17/2012
0.5ML	QCS ICV B	CPI	11C174-28549	9/17/2012
Prepared in 50ml 1%HNO3/5%HCl				

SDM 11/11/11

M3015

Metals Digestion Worksheet

Method Name 3015 Digestion **Prep Method** M3015 **Set** 111110A **Units** mL

Spikes	
Spiked ID 1	LCSW LOT# 1028408-29435
Spiked ID 2	LCSW LOT# 1028416-29433
Spiked ID 3	
Spiked ID 4	
Spiked By	NM Date: 11/10/11 10:40:00 AM
Witnessed By	KWS Date: 11/10/11 10:40:00 AM

Starting Temp:	25 C
Ending Temp:	170 C
Temperature Type:	Microwave
Sufficient Vol for Matrix QC:	YES
End Date/Time	11/10/11 12:00

	Sample	Sample Container	Spike Amount	Spike ID	Digested Amount	Final Volume	Start Date/Time	Comments
1	111110A Blk				45mL	50mL	11/10/11 10:40	equip: Venus
2	111110A LCS		90uL	1+2	45mL	50mL	11/10/11 10:40	equip: Venus
3	AY48273	AY48273W01			45mL	50mL	11/10/11 10:40	equip: Venus
4	AY48273 DUP	AY48273W01			45mL	50mL	11/10/11 10:40	equip: Venus
5	AY48273 MS	AY48273W01	90uL	1+2	45mL	50mL	11/10/11 10:40	equip: Venus
6	AY48639	AY48639W05			45mL	50mL	11/10/11 10:40	equip: Venus
7	AY48640	AY48640W05			45mL	50mL	11/10/11 10:40	equip: Venus
8	AY48641	AY48641W05			45mL	50mL	11/10/11 10:40	equip: Venus
9	AY48642	AY48642W05			45mL	50mL	11/10/11 10:40	equip: Venus
10	AY48643	AY48643W05			45mL	50mL	11/10/11 10:40	equip: Venus
11	AY48644	AY48644W02			45mL	50mL	11/10/11 10:40	equip: Venus
12	AY49333	AY49333W13			45mL	50mL	11/10/11 10:40	equip: Venus
13	AY49334	AY49334W51			45mL	50mL	11/10/11 10:40	equip: Venus
14	AY49334 MS	AY49334W52	90uL	1+2	45mL	50mL	11/10/11 10:40	equip: Venus
15	AY49334 MSD	AY49334W52	90uL	1+2	45mL	50mL	11/10/11 10:40	equip: Venus
16	AY49336	AY49336W13			45mL	50mL	11/10/11 10:40	equip: Venus
17	AY49481	AY49481W13			45mL	50mL	11/10/11 10:40	equip: Venus
18	AY49482	AY49482W13			45mL	50mL	11/10/11 10:40	equip: Venus
19	AY49559	AY49559W31			45mL	50mL	11/10/11 10:40	equip: Venus
20	AY49559 MS	AY49559W31	90uL	1+2	45mL	50mL	11/10/11 10:40	equip: Venus
21	AY49559 MSD	AY49559W31	90uL	1+2	45mL	50mL	11/10/11 10:40	equip: Venus
22	AY49561	AY49561W08			45mL	50mL	11/10/11 10:40	equip: Venus
23	AY49562	AY49562W08			45mL	50mL	11/10/11 10:40	equip: Venus

Solvent and Lot#
HNO3 J.T.B k19023 0095

Sample GOO Transfer	
Sample prep employee Initials	nm
Analyst's initials	NBS
Date	11-10-11
Time	13:00
Moved to	MSTAGS

Technicians Initials	
Scanned By	nm
Sample Preparation	lo
Digestion	lo
Bring up to volume	nm
Modified	11/10/11 10:19:53 AM

Reviewed By: [signature] **Date:** 11-10-11

M3015

Metals Digestion Worksheet

Method Name 3015 Digestion **Prep Method** M3015 **Set** 111110A **Units** mL

Spikes	
Spiked ID 1	LCSW LOT# 1028408-29435
Spiked ID 2	LCSW LOT# 1028416-29433
Spiked ID 3	
Spiked ID 4	
Spiked By	NM Date: 11/10/11 10:40:00 AM
Witnessed By	KWS Date: 11/10/11 10:40:00 AM

Starting Temp:	25 C
Ending Temp:	170 C
Temperature Type:	Microwave
Sufficient Vol for Matrix QC:	YES
End Date/Time	11/10/11 12:00

	Sample	Sample Container	Spike Amount	Spike ID	Digested Amount	Final Volume	Start Date/Time	Comments
24	AY50005	AY50005W08			45mL	50mL	11/10/11 10:40	equip: Venus

Solvent and Lot#	
HNO3 J.T.B k19023 0095	

Sample COC Transfer	
Sample prep employee Initials	nm
Analyst's Initials	NBS
Date	11-10-11
Time	13:00
Moved to	MSTACS

Technician's Initials	
Scanned By	nm
Sample Preparation	lo
Digestion	lo
Bring up to volume	nm
Modified	11/10/11 10:19:53 AM

Reviewed By: SA **Date:** 11-10-11

6020/200.8 Injection Log

Directory: K:\ICP-MS Optimus\raw data output csv\

RunID	Injected		Sample Name	Misc Info	FileName	Multiplier
1	11 Nov 2011	12:08	Calibration Blank		111111A	1.
2	11 Nov 2011	12:14	111111 Standard 1		111111A	1.
3	11 Nov 2011	12:20	111111 Standard 2		111111A	1.
4	11 Nov 2011	12:27	111111 Standard 3		111111A	1.
5	11 Nov 2011	12:33	111111 Standard 4		111111A	1.
6	11 Nov 2011	12:39	ICV 111111		111111A	1.
8	11 Nov 2011	12:57	ICB 111111		111111A	1.
9	11 Nov 2011	13:03	CCV 111111		111111A	1.
10	11 Nov 2011	13:09	CCB 111111		111111A	1.
11	11 Nov 2011	13:15	ICSA 111111		111111A	1.
12	11 Nov 2011	13:21	ICSAB 111111		111111A	1.
13	11 Nov 2011	13:33	CCV 111111		111111A	1.
14	11 Nov 2011	13:46	CCB 111111		111111A	1.
15	11 Nov 2011	14:16	111110A-3015-BLK		111111A	1.
16	11 Nov 2011	14:22	111110A-3015-LCS		111111A	1.
23	11 Nov 2011	15:05	CCV 111111		111111A	1.
24	11 Nov 2011	15:17	CCB 111111		111111A	1.
31	11 Nov 2011	16:00	AY49333W13		111111A	1.
32	11 Nov 2011	16:06	AY49334W51		111111A	1.
33	11 Nov 2011	16:12	AY49334W52 MS		111111A	1.
36	11 Nov 2011	16:30	CCV 111111		111111A	1.
37	11 Nov 2011	16:42	CCB 111111		111111A	1.
38	11 Nov 2011	16:48	AY49334W52 MSD		111111A	1.
39	11 Nov 2011	16:54	AY49334W51-A		111111A	1.
40	11 Nov 2011	17:00	AY49334W51-1/5		111111A	5.
49	11 Nov 2011	18:04	CCV 111111		111111A	1.
50	11 Nov 2011	18:16	CCB 111111		111111A	1.