



GC/MS Volatiles Case Narrative

Colorado Oil & Gas Conservation Commission

Complaint 200272771

Work Order Number: 1010128


1. This report consists of 2 water samples. The samples were received cool and intact by ALS on 10/08/10. All aqueous samples were free of headspace prior to analysis.
2. These samples were prepared according to SW-846, 3rd Edition procedures. Specifically, the water samples were prepared by purging 10 mL using purge and trap procedures based on Method 5030C.
3. The samples were analyzed using GC/MS with an RTX-624, RTX-VMS, or equivalent capillary column according to SOP 525 Revision 14 based on SW-846 Method 8260. All positive results were quantitated against the initial calibration standards using the internal standard technique. The identification of positive results was achieved by a comparison of the retention time and mass spectrum of the sample versus the daily calibration standard.
4. All initial calibration criteria for SPCC's and CCC's were met. If average response factors were used in the initial calibration, %RSD was $\leq 15\%$. If linear or higher order regression calibrations were used in the initial calibration, the coefficient of determination (r^2) ≥ 0.99 .
5. All initial calibrations are verified by comparing a second source standard calibration verification (ICV) against the calibration curve. All target compounds in the second source verification had a %D of less than 25%.
6. All criteria for SPCC's and CCC's were met in daily (continuing) calibration verifications (CCV).
7. Methylene chloride, acetone and 2-butanone are common laboratory contaminants. In order to minimize the levels of these compounds detected in the gc/ms analysis, ALS has designated its volatile laboratory as a restricted access area. In addition, the laboratory has been equipped with a dedicated, air intake and exhaust system that operates under positive pressure in order to minimize cross contamination of these compounds. Due to fluctuations in ambient laboratory conditions, reported sample values for common laboratory contaminants may be due to lab contamination even if the compound in question is not detected in the associated method blank.



All method blank criteria were met.

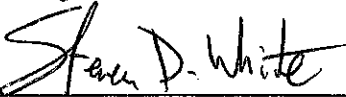
8. All laboratory control sample and laboratory control sample duplicate recoveries and RPDs were within the acceptance criteria.
9. Since a sample from this order number was not the selected quality control (QC) sample, matrix specific QC results are not included in this report.
10. The samples were analyzed within the established holding time.
11. All surrogate recoveries were within acceptance criteria.
12. All internal standard recoveries were within acceptance criteria.
13. Manual integrations are performed when needed to provide consistent and defensible data following the guidelines in SOP 939 Revision 3.

The data contained in the following report have been reviewed and approved by the personnel listed below. In addition, ALS certifies that the analyses reported herein are true, complete and correct within the limits of the methods employed.



Sharon L. Jobs
Organics Primary Data Reviewer

10-26-10
Date



Stan D. White
Organics Final Data Reviewer

10/26/10
Date



ALS
Data Qualifier Flags
Chromatography and Mass Spectrometry

- U or ND:** This flag indicates that the compound was analyzed for but not detected.
- J:** This flag indicates an estimated value. This flag is used as follows : (1) when estimating a concentration for tentatively identified compounds (TICs) where a 1:1 response is assumed; (2) when the mass spectral and retention time data indicate the presence of a compound that meets the volatile and semivolatile GC/MS identification criteria, and the result is less than the reporting limit (RL) but greater than the method detection limit (MDL); (3) when the retention time data indicate the presence of a compound that meets the GC identification criteria, and the result is less than the RL but greater than the MDL; and (4) the reported value is estimated.
- B:** This flag is used when the analyte is detected in the associated method blank as well as in the sample. It indicates probable blank contamination and warns the data user. This flag shall be used for a tentatively identified compound (TIC) as well as for a positively identified target compound.
- E:** This flag identifies compounds whose concentration exceeds the upper level of the calibration range.
- A:** This flag indicates that a tentatively identified compound is a suspected aldol-condensation product.
- X:** This flag indicates that the analyte was diluted below an accurate quantitation level.
- *:** This flag indicates that a spike recovery is equal to or outside the control criteria used.
- +:** This flag indicates that the relative percent difference (RPD) equals or exceeds the control criteria.

ALS Environmental -- FC

Sample Number(s) Cross-Reference Table

Paragon OrderNum: 1010128

Client Name: Colorado Oil & Gas Conservation Commission

Client Project Name: Complaint 200272771

Client Project Number:

Client PO Number: OE PHA 11000000014

Client Sample Number	Lab Sample Number	COC Number	Matrix	Date Collected	Time Collected
Burge WW	1010128-1		WATER	06-Oct-10	11:25
Trip Blank	1010128-2		WATER	06-Oct-10	



ALS Laboratory Group

225 Commerce Drive, Fort Collins, Colorado 80524
TF: (800) 443-1511 PH: (970) 490-1511 FX: (970) 490-1522

Chain-of-Custody

Form 202r8

WORKORDER #

1010128

PAGE

of 1

DISPOSAL

By Lab or Return to Client

PROJECT NAME Complant 20027271

SAMPLER

SITE ID

DATE

TURNAROUND

14 day

PROJECT No.

EDD FORMAT

PURCHASE ORDER

PHA 100611-14

COMPANY NAME Colo Oil & Gas Cns. Lm

BILL TO COMPANY

SEND REPORT TO Peter Gintantes

INVOICE ATTN TO

ADDRESS PO Box 108

ADDRESS

CITY / STATE / ZIP Trinidad CO 81082

CITY / STATE / ZIP

PHONE 719-846-3091

PHONE

FAX

FAX

E-MAIL peter.gintantes@state.co.us

E-MAIL

Lab ID

Field ID

Matrix

Sample Date

Sample Time

Bottles

Pres.

QC

1 Burge WW
Burge WW

W

10/6/10 11:25A

1

X X X X X X X X X X

2 Trip Blank

W

1

1

X

Anions = Br, Cl, F, NO₂, NO₃, SO₄
200.8 = Al, Sb, As, Cd, Pb, Mo, Se, Ag, Te, U
200.7 = Ba, Be, B, Ca, Cr, Co, Cu, Fe, Li, Mg, Mn, Ni, K, Si, Na, Sr, Zn

Time Zone (Circle): EST CST PST Matrix: O = oil S = soil NS = non-soil solid W = water L = liquid E = extract F = filter

For metals or anions, please detail analytes below.

Comments:

Dissolved = Filter + preserve at lab.

QC PACKAGE (check below)

☒ LEVEL II (Standard QC)
☐ LEVEL III (Std QC + forms)
☐ LEVEL IV (Std QC + forms + raw data)

Preservative Key: 1-HCl 2-HNO₃ 3-H₂SO₄ 4-NaOH 5-NaHSO₄ 7-Other 8-4 degrees C 9-5035

SIGNATURE

PRINTED NAME

DATE

TIME

RELINQUISHED BY

RECEIVED BY

RELINQUISHED BY

RECEIVED BY

RELINQUISHED BY

RECEIVED BY

Angela Bellumani

C. Cochran

10/7/10 9:00A

10/8/10 0945



CONDITION OF SAMPLE UPON RECEIPT FORM

Client: COGCCWorkorder No: 1010128Project Manager: AWInitials: CW Date: 10-8-10

1. Does this project require any special handling in addition to standard Paragon procedures?		<u>YES</u>	NO
2. Are custody seals on shipping containers intact?	NONE	<u>YES</u>	NO
3. Are Custody seals on sample containers intact?	NONE	<u>YES</u>	NO
4. Is there a COC (Chain-of-Custody) present or other representative documents?		<u>YES</u>	NO
5. Are the COC and bottle labels complete and legible?		<u>YES</u>	NO
6. Is the COC in agreement with samples received? (IDs, dates, times, no. of samples, no. of containers, matrix, requested analyses, etc.)		<u>YES</u>	NO
7. Were airbills / shipping documents present and/or removable?	DROP OFF	<u>YES</u>	NO
8. Are all aqueous samples requiring preservation preserved correctly? (excluding volatiles)	N/A	<u>YES</u>	NO
9. Are all aqueous non-preserved samples pH 4-9?	N/A	<u>YES</u>	NO
10. Is there sufficient sample for the requested analyses?		<u>YES</u>	NO
11. Were all samples placed in the proper containers for the requested analyses?		<u>YES</u>	NO
12. Are all samples within holding times for the requested analyses?		<u>YES</u>	NO
13. Were all sample containers received intact ? (not broken or leaking, etc.)		<u>YES</u>	NO
14. Are all samples requiring no headspace (VOC, GRO, RSK/MEE, Rx CN/S, radon) headspace free? Size of bubble: <u> </u> < green pea <u> </u> > green pea	N/A	<u>YES</u>	NO
15. Do perchlorate LCMS-MS samples have headspace ? (at least 1/3 of container required)	N/A	<u>YES</u>	NO
16. Were samples checked for and free from the presence of residual chlorine ? (Applicable when PM has indicated samples are from a chlorinated water source; note if field preservation with sodium thiosulfate was not observed.)	N/A	<u>YES</u>	NO
17. Were the samples shipped on ice ?		<u>YES</u>	NO
18. Were cooler temperatures measured at 0.1-6.0°C? IR gun used*: <u>#2</u> <u>#4</u>	RAD ONLY	<u>YES</u>	NO
Cooler #: <u>1</u>			
Temperature (°C): <u>4.5</u>			
No. of custody seals on cooler: <u>1</u>			
DOT Survey/ Acceptance Information	External µR/hr reading: <u>12</u>		
	Background µR/hr reading: <u>11</u>		
Were external µR/hr readings ≤ two times background and within DOT acceptance criteria? <u>YES</u> / NO / NA (If no, see Form 008.)			

Additional Information: PROVIDE DETAILS BELOW FOR A NO RESPONSE TO ANY QUESTION ABOVE, EXCEPT #1 AND #16

received out of hotel with only a few hours of hold time remaining.
AW 10/8/10

If applicable, was the client contacted? YES / NO / NA Contact: _____ Date/Time: _____

Project Manager Signature / Date: _____

GC/MS Volatiles

Method SW8260_2514

Method Blank

Lab Name: ALS Environmental -- FC

Work Order Number: 1010128

Client Name: Colorado Oil & Gas Conservation Commission

ClientProject ID: Complaint 200272771

Lab ID: VL101011-2MB

Sample Matrix: WATER

% Moisture: N/A

Date Collected: N/A

Date Extracted: 11-Oct-10

Date Analyzed: 11-Oct-10

Prep Method: SW5030 Rev C

Prep Batch: VL101011-2

QCBatchID: VL101011-2-3

Run ID: VL101011-2A

Cleanup: NONE

Basis: N/A

File Name: B68264

Sample Aliquot: 10 ml

Final Volume: 10 ml

Result Units: UG/L

Clean DF: 1

CASNO	Target Analyte	DF	Result	Reporting Limit	Result Qualifier	EPA Qualifier
75-71-8	DICHLORODIFLUOROMETHANE	1	1	1	U	
74-87-3	CHLOROMETHANE	1	1	1	U	
75-01-4	VINYL CHLORIDE	1	1	1	U	
74-83-9	BROMOMETHANE	1	1	1	U	
75-00-3	CHLOROETHANE	1	1	1	U	
75-69-4	TRICHLOROFLUOROMETHANE	1	1	1	U	
75-35-4	1,1-DICHLOROETHENE	1	1	1	U	
76-13-1	1,1,2-TRICHLORO-1,2,2-TRIFLUOROET	1	1	1	U	
67-64-1	ACETONE	1	10	10	U	
74-88-4	IODOMETHANE	1	1	1	U	
75-15-0	CARBON DISULFIDE	1	1	1	U	
75-09-2	METHYLENE CHLORIDE	1	1	1	U	
156-60-5	TRANS-1,2-DICHLOROETHENE	1	1	1	U	
1634-04-4	METHYL TERTIARY BUTYL ETHER	1	1	1	U	
75-34-3	1,1-DICHLOROETHANE	1	1	1	U	
108-05-4	VINYL ACETATE	1	2	2	U	
156-59-2	CIS-1,2-DICHLOROETHENE	1	1	1	U	
78-93-3	2-BUTANONE	1	10	10	U	
74-97-5	BROMOCHLOROMETHANE	1	1	1	U	
67-66-3	CHLOROFORM	1	1	1	U	
71-55-6	1,1,1-TRICHLOROETHANE	1	1	1	U	
594-20-7	2,2-DICHLOROPROPANE	1	1	1	U	
56-23-5	CARBON TETRACHLORIDE	1	1	1	U	
563-58-6	1,1-DICHLOROPROPENE	1	1	1	U	
107-06-2	1,2-DICHLOROETHANE	1	1	1	U	
71-43-2	BENZENE	1	1	1	U	
79-01-6	TRICHLOROETHENE	1	1	1	U	

Data Package ID: VL1010128-1

Date Printed: Tuesday, October 26, 2010

ALS Environmental -- FC

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LIMS Version: 6.423A

GC/MS Volatiles

Method SW8260_2514

Method Blank

Lab Name: ALS Environmental -- FC

Work Order Number: 1010128

Client Name: Colorado Oil & Gas Conservation Commission

ClientProject ID: Complaint 200272771

Lab ID: VL101011-2MB

Sample Matrix: WATER

% Moisture: N/A

Date Collected: N/A

Date Extracted: 11-Oct-10

Date Analyzed: 11-Oct-10

Prep Method: SW5030 Rev C

Prep Batch: VL101011-2

QCBatchID: VL101011-2-3

Run ID: VL101011-2A

Cleanup: NONE

Basis: N/A

File Name: B68264

Sample Aliquot: 10 ml

Final Volume: 10 ml

Result Units: UG/L

Clean DF: 1

78-87-5	1,2-DICHLOROPROPANE	1	1	1	U	
74-95-3	DIBROMOMETHANE	1	1	1	U	
75-27-4	BROMODICHLOROMETHANE	1	1	1	U	
10061-01-5	CIS-1,3-DICHLOROPROPENE	1	1	1	U	
108-10-1	4-METHYL-2-PENTANONE	1	10	10	U	
108-88-3	TOLUENE	1	1	1	U	
10061-02-6	TRANS-1,3-DICHLOROPROPENE	1	1	1	U	
79-00-5	1,1,2-TRICHLOROETHANE	1	1	1	U	
591-78-6	2-HEXANONE	1	10	10	U	
127-18-4	TETRACHLOROETHENE	1	1	1	U	
142-28-9	1,3-DICHLOROPROPANE	1	1	1	U	
124-48-1	DIBROMOCHLOROMETHANE	1	1	1	U	
106-93-4	1,2-DIBROMOETHANE	1	1	1	U	
544-10-5	1-CHLOROHEXANE	1	1	1	U	
108-90-7	CHLOROBENZENE	1	1	1	U	
630-20-6	1,1,1,2-TETRACHLOROETHANE	1	1	1	U	
100-41-4	ETHYLBENZENE	1	1	1	U	
136777-61-2	M+P-XYLENE	1	1	1	U	
95-47-6	O-XYLENE	1	1	1	U	
100-42-5	STYRENE	1	1	1	U	
75-25-2	BROMOFORM	1	1	1	U	
98-82-8	ISOPROPYLBENZENE	1	1	1	U	
96-18-4	1,2,3-TRICHLOROPROPANE	1	1	1	U	
79-34-5	1,1,2,2-TETRACHLOROETHANE	1	1	1	U	
108-86-1	BROMOBENZENE	1	1	1	U	
103-65-1	N-PROPYLBENZENE	1	1	1	U	
95-49-8	2-CHLOROTOLUENE	1	1	1	U	
108-67-8	1,3,5-TRIMETHYLBENZENE	1	1	1	U	
106-43-4	4-CHLOROTOLUENE	1	1	1	U	

Data Package ID: VL1010128-1

Date Printed: Tuesday, October 26, 2010

ALS Environmental -- FC

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GC/MS Volatiles

Method SW8260_2514

Method Blank

Lab Name: ALS Environmental -- FC

Work Order Number: 1010128

Client Name: Colorado Oil & Gas Conservation Commission

ClientProject ID: Complaint 200272771

Lab ID: VL101011-2MB

Sample Matrix: WATER

% Moisture: N/A

Date Collected: N/A

Date Extracted: 11-Oct-10

Date Analyzed: 11-Oct-10

Prep Method: SW5030 Rev C

Prep Batch: VL101011-2

QCBatchID: VL101011-2-3

Run ID: VL101011-2A

Cleanup: NONE

Basis: N/A

File Name: B68264

Sample Aliquot: 10 ml

Final Volume: 10 ml

Result Units: UG/L

Clean DF: 1

98-06-6	TERT-BUTYLBENZENE	1	1	1	U	
95-63-6	1,2,4-TRIMETHYLBENZENE	1	1	1	U	
135-98-8	SEC-BUTYLBENZENE	1	1	1	U	
541-73-1	1,3-DICHLOROBENZENE	1	1	1	U	
99-87-6	P-ISOPROPYLTOLUENE	1	1	1	U	
106-46-7	1,4-DICHLOROBENZENE	1	1	1	U	
104-51-8	N-BUTYLBENZENE	1	1	1	U	
95-50-1	1,2-DICHLOROBENZENE	1	1	1	U	
96-12-8	1,2-DIBROMO-3-CHLOROPROPANE	1	2	2	U	
120-82-1	1,2,4-TRICHLOROBENZENE	1	1	1	U	
87-68-3	HEXACHLOROBUTADIENE	1	1	1	U	
91-20-3	NAPHTHALENE	1	1	1	U	
87-61-6	1,2,3-TRICHLOROBENZENE	1	1	1	U	

Surrogate Recovery

CASNO	Surrogate Analyte	Result	Flag	Spike Amount	Percent Recovery	Control Limits
460-00-4	4-BROMOFLUOROBENZENE	23.3		25	93	87 - 108
1868-53-7	DIBROMOFLUOROMETHANE	23.8		25	95	89 - 112
2037-26-5	TOLUENE-D8	25.5		25	102	87 - 109

Data Package ID: VL1010128-1

Date Printed: Tuesday, October 26, 2010

ALS Environmental -- FC

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GC/MS Volatiles

Method SW8260_25

Tentatively Identified Compounds

Lab Name: ALS Environmental -- FC

Work Order Number: 1010128

Client Name: Colorado Oil & Gas Conservation Commission

ClientProject ID: Complaint 200272771

Field ID:	
Lab ID:	VL101011-2MB

Sample Matrix: WATER

% Moisture: N/A

Date Collected: N/A

Date Extracted: 11-Oct-10

Date Analyzed: 11-Oct-10

Prep Batch: VL101011-2

QCBatchID: VL101011-2-3

Run ID: VL101011-2A

Cleanup: NONE

Basis: As Received

Sample Aliquot: 10 ml

Final Volume: 10 ml

Clean DF: 1

File Name: B68264

CASNO	Retention Time	Target Analyte	Dilution Factor	Result	Units	Qualifier
		NONE DETECTED	1			U

Data Package ID: VL1010128-1

GC/MS Volatiles

Method SW8260_2514

Sample Results

Lab Name: ALS Environmental -- FC

Work Order Number: 1010128

Client Name: Colorado Oil & Gas Conservation Commission

ClientProject ID: Complaint 200272771

Field ID: Burge WW

Lab ID: 1010128-1

Sample Matrix: WATER

% Moisture: N/A

Date Collected: 06-Oct-10

Date Extracted: 11-Oct-10

Date Analyzed: 11-Oct-10

Prep Method: SW5030 Rev C

Prep Batch: VL101011-2

QCBatchID: VL101011-2-3

Run ID: VL101011-2A

Cleanup: NONE

Basis: As Received

File Name: B68278

Sample Aliquot: 10 ml

Final Volume: 10 ml

Result Units: UG/L

Clean DF: 1

CASNO	Target Analyte	Dilution Factor	Result	Reporting Limit	Result Qualifier	EPA Qualifier
75-71-8	DICHLORODIFLUOROMETHANE	1	1	1	U	
74-87-3	CHLOROMETHANE	1	1	1	U	
75-01-4	VINYL CHLORIDE	1	1	1	U	
74-83-9	BROMOMETHANE	1	1	1	U	
75-00-3	CHLOROETHANE	1	1	1	U	
75-69-4	TRICHLOROFLUOROMETHANE	1	1	1	U	
75-35-4	1,1-DICHLOROETHENE	1	1	1	U	
76-13-1	1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE	1	1	1	U	
67-64-1	ACETONE	1	10	10	U	
74-88-4	IODOMETHANE	1	1	1	U	
75-15-0	CARBON DISULFIDE	1	1	1	U	
75-09-2	METHYLENE CHLORIDE	1	1	1	U	
156-60-5	TRANS-1,2-DICHLOROETHENE	1	1	1	U	
1634-04-4	METHYL TERTIARY BUTYL ETHER	1	1	1	U	
75-34-3	1,1-DICHLOROETHANE	1	1	1	U	
108-05-4	VINYL ACETATE	1	2	2	U	
156-59-2	CIS-1,2-DICHLOROETHENE	1	1	1	U	
78-93-3	2-BUTANONE	1	10	10	U	
74-97-5	BROMOCHLOROMETHANE	1	1	1	U	
67-66-3	CHLOROFORM	1	1	1	U	
71-55-6	1,1,1-TRICHLOROETHANE	1	1	1	U	
594-20-7	2,2-DICHLOROPROPANE	1	1	1	U	
56-23-5	CARBON TETRACHLORIDE	1	1	1	U	
563-58-6	1,1-DICHLOROPROPENE	1	1	1	U	
107-06-2	1,2-DICHLOROETHANE	1	1	1	U	
71-43-2	BENZENE	1	1	1	U	
79-01-6	TRICHLOROETHENE	1	1	1	U	

Data Package ID: VL1010128-1

Date Printed: Tuesday, October 26, 2010

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LIMS Version: 6.423A

GC/MS Volatiles

Method SW8260_2514

Sample Results

Lab Name: ALS Environmental -- FC

Work Order Number: 1010128

Client Name: Colorado Oil & Gas Conservation Commission

ClientProject ID: Complaint 200272771

Field ID: Burge WW

Lab ID: 1010128-1

Sample Matrix: WATER

% Moisture: N/A

Date Collected: 06-Oct-10

Date Extracted: 11-Oct-10

Date Analyzed: 11-Oct-10

Prep Method: SW5030 Rev C

Prep Batch: VL101011-2

QCBatchID: VL101011-2-3

Run ID: VL101011-2A

Cleanup: NONE

Basis: As Received

File Name: B68278

Sample Aliquot: 10 ml

Final Volume: 10 ml

Result Units: UG/L

Clean DF: 1

78-87-5	1,2-DICHLOROPROPANE	1	1	1	U	
74-95-3	DIBROMOMETHANE	1	1	1	U	
75-27-4	BROMODICHLOROMETHANE	1	1	1	U	
10061-01-5	CIS-1,3-DICHLOROPROPENE	1	1	1	U	
108-10-1	4-METHYL-2-PENTANONE	1	10	10	U	
108-88-3	TOLUENE	1	1	1	U	
10061-02-6	TRANS-1,3-DICHLOROPROPENE	1	1	1	U	
79-00-5	1,1,2-TRICHLOROETHANE	1	1	1	U	
591-78-6	2-HEXANONE	1	10	10	U	
127-18-4	TETRACHLOROETHENE	1	1	1	U	
142-28-9	1,3-DICHLOROPROPANE	1	1	1	U	
124-48-1	DIBROMOCHLOROMETHANE	1	1	1	U	
106-93-4	1,2-DIBROMOETHANE	1	1	1	U	
544-10-5	1-CHLOROHEXANE	1	1	1	U	
108-90-7	CHLOROBENZENE	1	1	1	U	
630-20-6	1,1,1,2-TETRACHLOROETHANE	1	1	1	U	
100-41-4	ETHYLBENZENE	1	1	1	U	
136777-61-2	M+P-XYLENE	1	1	1	U	
95-47-6	O-XYLENE	1	1	1	U	
100-42-5	STYRENE	1	1	1	U	
75-25-2	BROMOFORM	1	1	1	U	
98-82-8	ISOPROPYLBENZENE	1	1	1	U	
96-18-4	1,2,3-TRICHLOROPROPANE	1	1	1	U	
79-34-5	1,1,2,2-TETRACHLOROETHANE	1	1	1	U	
108-86-1	BROMOBENZENE	1	1	1	U	
103-65-1	N-PROPYLBENZENE	1	1	1	U	
95-49-8	2-CHLOROTOLUENE	1	1	1	U	
108-67-8	1,3,5-TRIMETHYLBENZENE	1	1	1	U	
106-43-4	4-CHLOROTOLUENE	1	1	1	U	

Data Package ID: VL1010128-1

Date Printed: Tuesday, October 26, 2010

ALS Environmental -- FC

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LIMS Version: 6.423A

GC/MS Volatiles

Method SW8260_2514

Sample Results

Lab Name: ALS Environmental -- FC

Work Order Number: 1010128

Client Name: Colorado Oil & Gas Conservation Commission

ClientProject ID: Complaint 200272771

Field ID:	Burge WW
Lab ID:	1010128-1

Sample Matrix: WATER
% Moisture: N/A
Date Collected: 06-Oct-10
Date Extracted: 11-Oct-10
Date Analyzed: 11-Oct-10
Prep Method: SW5030 Rev C

Prep Batch: VL101011-2
QCBatchID: VL101011-2-3
Run ID: VL101011-2A
Cleanup: NONE
Basis: As Received
File Name: B68278

Sample Aliquot: 10 ml
Final Volume: 10 ml
Result Units: UG/L
Clean DF: 1

98-06-6	TERT-BUTYLBENZENE	1	1	1	U	
95-63-6	1,2,4-TRIMETHYLBENZENE	1	1	1	U	
135-98-8	SEC-BUTYLBENZENE	1	1	1	U	
541-73-1	1,3-DICHLOROBENZENE	1	1	1	U	
99-87-6	P-ISOPROPYLTOLUENE	1	1	1	U	
106-46-7	1,4-DICHLOROBENZENE	1	1	1	U	
104-51-8	N-BUTYLBENZENE	1	1	1	U	
95-50-1	1,2-DICHLOROBENZENE	1	1	1	U	
96-12-8	1,2-DIBROMO-3-CHLOROPROPANE	1	2	2	U	
120-82-1	1,2,4-TRICHLOROBENZENE	1	1	1	U	
87-68-3	HEXACHLOROBUTADIENE	1	1	1	U	
91-20-3	NAPHTHALENE	1	1	1	U	
87-61-6	1,2,3-TRICHLOROBENZENE	1	1	1	U	

Surrogate Recovery

CASNO	Surrogate Analyte	Result	Flag	Spike Amount	Percent Recovery	Control Limits
460-00-4	4-BROMOFLUOROBENZENE	24		25	96	87 - 108
1868-53-7	DIBROMOFLUOROMETHANE	24.8		25	99	89 - 112
2037-26-5	TOLUENE-D8	24.6		25	98	87 - 109

Data Package ID: VL1010128-1

Date Printed: Tuesday, October 26, 2010

ALS Environmental -- FC

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LIMS Version: 6.423A

GC/MS Volatiles

Method SW8260_25

Tentatively Identified Compounds

Lab Name: ALS Environmental -- FC

Work Order Number: 1010128

Client Name: Colorado Oil & Gas Conservation Commission

ClientProject ID: Complaint 200272771

Field ID: Burge WW

Lab ID: 1010128-1

Sample Matrix: WATER

% Moisture: N/A

Date Collected: 06-Oct-10

Date Extracted: 11-Oct-10

Date Analyzed: 11-Oct-10

Prep Batch: VL101011-2

QCBatchID: VL101011-2-3

Run ID: VL101011-2A

Cleanup: NONE

Basis: As Received

Sample Aliquot: 10 ml

Final Volume: 10 ml

Clean DF: 1

File Name: B68278

CASNO	Retention Time	Target Analyte	Dilution Factor	Result	Units	Qualifier
109-99-9	8.51	TETRAHYDROFURAN1	1	0.67	UG/L	J

Data Package ID: VL1010128-1

GC/MS Volatiles

Method SW8260_2514

Sample Results

Lab Name: ALS Environmental -- FC

Work Order Number: 1010128

Client Name: Colorado Oil & Gas Conservation Commission

ClientProject ID: Complaint 200272771

Field ID: Trip Blank

Lab ID: 1010128-2

Sample Matrix: WATER

% Moisture: N/A

Date Collected: 06-Oct-10

Date Extracted: 11-Oct-10

Date Analyzed: 11-Oct-10

Prep Method: SW5030 Rev C

Prep Batch: VL101011-2

QCBatchID: VL101011-2-3

Run ID: VL101011-2A

Cleanup: NONE

Basis: As Received

File Name: B68279

Sample Aliquot: 10 ml

Final Volume: 10 ml

Result Units: UG/L

Clean DF: 1

CASNO	Target Analyte	Dilution Factor	Result	Reporting Limit	Result Qualifier	EPA Qualifier
75-71-8	DICHLORODIFLUOROMETHANE	1	1	1	U	
74-87-3	CHLOROMETHANE	1	1	1	U	
75-01-4	VINYL CHLORIDE	1	1	1	U	
74-83-9	BROMOMETHANE	1	1	1	U	
75-00-3	CHLOROETHANE	1	1	1	U	
75-69-4	TRICHLOROFLUOROMETHANE	1	1	1	U	
75-35-4	1,1-DICHLOROETHENE	1	1	1	U	
76-13-1	1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHANE	1	1	1	U	
67-64-1	ACETONE	1	10	10	U	
74-88-4	IODOMETHANE	1	1	1	U	
75-15-0	CARBON DISULFIDE	1	1	1	U	
75-09-2	METHYLENE CHLORIDE	1	1	1	U	
156-60-5	TRANS-1,2-DICHLOROETHENE	1	1	1	U	
1634-04-4	METHYL TERTIARY BUTYL ETHER	1	1	1	U	
75-34-3	1,1-DICHLOROETHANE	1	1	1	U	
108-05-4	VINYL ACETATE	1	2	2	U	
156-59-2	CIS-1,2-DICHLOROETHENE	1	1	1	U	
78-93-3	2-BUTANONE	1	10	10	U	
74-97-5	BROMOCHLOROMETHANE	1	1	1	U	
67-66-3	CHLOROFORM	1	1	1	U	
71-55-6	1,1,1-TRICHLOROETHANE	1	1	1	U	
594-20-7	2,2-DICHLOROPROPANE	1	1	1	U	
56-23-5	CARBON TETRACHLORIDE	1	1	1	U	
563-58-6	1,1-DICHLOROPROPENE	1	1	1	U	
107-06-2	1,2-DICHLOROETHANE	1	1	1	U	
71-43-2	BENZENE	1	1	1	U	
79-01-6	TRICHLOROETHENE	1	1	1	U	

Data Package ID: VL1010128-1

Date Printed: Tuesday, October 26, 2010

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GC/MS Volatiles

Method SW8260_2514

Sample Results

Lab Name: ALS Environmental -- FC

Work Order Number: 1010128

Client Name: Colorado Oil & Gas Conservation Commission

ClientProject ID: Complaint 200272771

Field ID:	Trip Blank
Lab ID:	1010128-2

Sample Matrix: WATER
% Moisture: N/A
Date Collected: 06-Oct-10
Date Extracted: 11-Oct-10
Date Analyzed: 11-Oct-10
Prep Method: SW5030 Rev C

Prep Batch: VL101011-2
QCBatchID: VL101011-2-3
Run ID: VL101011-2A
Cleanup: NONE
Basis: As Received
File Name: B68279

Sample Aliquot: 10 ml
Final Volume: 10 ml
Result Units: UG/L
Clean DF: 1

78-87-5	1,2-DICHLOROPROPANE	1	1	1	U	
74-95-3	DIBROMOMETHANE	1	1	1	U	
75-27-4	BROMODICHLOROMETHANE	1	1	1	U	
10061-01-5	CIS-1,3-DICHLOROPROPENE	1	1	1	U	
108-10-1	4-METHYL-2-PENTANONE	1	10	10	U	
108-88-3	TOLUENE	1	1	1	U	
10061-02-6	TRANS-1,3-DICHLOROPROPENE	1	1	1	U	
79-00-5	1,1,2-TRICHLOROETHANE	1	1	1	U	
591-78-6	2-HEXANONE	1	10	10	U	
127-18-4	TETRACHLOROETHENE	1	1	1	U	
142-28-9	1,3-DICHLOROPROPANE	1	1	1	U	
124-48-1	DIBROMOCHLOROMETHANE	1	1	1	U	
106-93-4	1,2-DIBROMOETHANE	1	1	1	U	
544-10-5	1-CHLOROHEXANE	1	1	1	U	
108-90-7	CHLOROBENZENE	1	1	1	U	
630-20-6	1,1,1,2-TETRACHLOROETHANE	1	1	1	U	
100-41-4	ETHYLBENZENE	1	1	1	U	
136777-61-2	M+P-XYLENE	1	1	1	U	
95-47-6	O-XYLENE	1	1	1	U	
100-42-5	STYRENE	1	1	1	U	
75-25-2	BROMOFORM	1	1	1	U	
98-82-8	ISOPROPYLBENZENE	1	1	1	U	
96-18-4	1,2,3-TRICHLOROPROPANE	1	1	1	U	
79-34-5	1,1,2,2-TETRACHLOROETHANE	1	1	1	U	
108-86-1	BROMOBENZENE	1	1	1	U	
103-65-1	N-PROPYLBENZENE	1	1	1	U	
95-49-8	2-CHLOROTOLUENE	1	1	1	U	
108-67-8	1,3,5-TRIMETHYLBENZENE	1	1	1	U	
106-43-4	4-CHLOROTOLUENE	1	1	1	U	

Data Package ID: VL1010128-1

GC/MS Volatiles

Method SW8260_2514

Sample Results

Lab Name: ALS Environmental -- FC

Work Order Number: 1010128

Client Name: Colorado Oil & Gas Conservation Commission

ClientProject ID: Complaint 200272771

Field ID:	Trip Blank
Lab ID:	1010128-2

Sample Matrix: WATER
% Moisture: N/A
Date Collected: 06-Oct-10
Date Extracted: 11-Oct-10
Date Analyzed: 11-Oct-10
Prep Method: SW5030 Rev C

Prep Batch: VL101011-2
QCBatchID: VL101011-2-3
Run ID: VL101011-2A
Cleanup: NONE
Basis: As Received
File Name: B68279

Sample Aliquot: 10 ml
Final Volume: 10 ml
Result Units: UG/L
Clean DF: 1

98-06-6	TERT-BUTYLBENZENE	1	1	1	U	
95-63-6	1,2,4-TRIMETHYLBENZENE	1	1	1	U	
135-98-8	SEC-BUTYLBENZENE	1	1	1	U	
541-73-1	1,3-DICHLOROBENZENE	1	1	1	U	
99-87-6	P-ISOPROPYLTOLUENE	1	1	1	U	
106-46-7	1,4-DICHLOROBENZENE	1	1	1	U	
104-51-8	N-BUTYLBENZENE	1	1	1	U	
95-50-1	1,2-DICHLOROBENZENE	1	1	1	U	
96-12-8	1,2-DIBROMO-3-CHLOROPROPANE	1	2	2	U	
120-82-1	1,2,4-TRICHLOROBENZENE	1	1	1	U	
87-68-3	HEXACHLOROBUTADIENE	1	1	1	U	
91-20-3	NAPHTHALENE	1	1	1	U	
87-61-6	1,2,3-TRICHLOROBENZENE	1	1	1	U	

Surrogate Recovery

CASNO	Surrogate Analyte	Result	Flag	Spike Amount	Percent Recovery	Control Limits
460-00-4	4-BROMOFLUOROBENZENE	24		25	96	87 - 108
1868-53-7	DIBROMOFLUOROMETHANE	23.4		25	94	89 - 112
2037-26-5	TOLUENE-D8	25		25	100	87 - 109

Data Package ID: VL1010128-1

Date Printed: Tuesday, October 26, 2010

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GC/MS Volatiles

Method SW8260_25

Tentatively Identified Compounds

Lab Name: ALS Environmental -- FC

Work Order Number: 1010128

Client Name: Colorado Oil & Gas Conservation Commission

ClientProject ID: Complaint 200272771

Field ID:	Trip Blank
Lab ID:	1010128-2

Sample Matrix: WATER

% Moisture: N/A

Date Collected: 06-Oct-10

Date Extracted: 11-Oct-10

Date Analyzed: 11-Oct-10

Prep Batch: VL101011-2

QCBatchID: VL101011-2-3

Run ID: VL101011-2A

Cleanup: NONE

Basis: As Received

Sample Aliquot: 10 ml

Final Volume: 10 ml

Clean DF: 1

File Name: B68279

CASNO	Retention Time	Target Analyte	Dilution Factor	Result	Units	Qualifier
		NONE DETECTED	1			U

Data Package ID: VL1010128-1

GC/MS Volatiles

Method SW8260_2514

Laboratory Control Sample and Laboratory Control Sample Duplicate

Lab Name: ALS Environmental -- FC

Work Order Number: 1010128

Client Name: Colorado Oil & Gas Conservation Commission

ClientProject ID: Complaint 200272771

Lab ID: VL101011-2LCS

Sample Matrix: WATER

% Moisture: N/A

Date Collected: N/A

Date Extracted: 10/11/2010

Date Analyzed: 10/11/2010

Prep Method: SW5030C

Prep Batch: VL101011-2

QCBatchID: VL101011-2-3

Run ID: VL101011-2A

Cleanup: NONE

Basis: N/A

File Name: B68261

Sample Aliquot: 10 ml

Final Volume: 10 ml

Result Units: UG/L

Clean DF: 1

CASNO	Target Analyte	Spike Added	LCS Result	Reporting Limit	Result Qualifier	LCS % Rec.	Control Limits
75-71-8	DICHLORODIFLUOROMETHANE	10	9.51	1		95	63 - 125%
74-87-3	CHLOROMETHANE	10	10.4	1		104	73 - 122%
75-01-4	VINYL CHLORIDE	10	11.9	1		119	72 - 123%
74-83-9	BROMOMETHANE	10	10.3	1		103	68 - 123%
75-00-3	CHLOROETHANE	10	9.55	1		95	74 - 124%
75-69-4	TRICHLOROFLUOROMETHANE	10	9.75	1		97	74 - 124%
75-35-4	1,1-DICHLOROETHENE	10	9.59	1		96	77 - 119%
76-13-1	1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHA	10	10.4	1		104	79 - 122%
67-64-1	ACETONE	40	31.4	10		79	62 - 142%
74-88-4	IODOMETHANE	10	9.98	1		100	72 - 126%
75-15-0	CARBON DISULFIDE	10	10.4	1		104	76 - 121%
75-09-2	METHYLENE CHLORIDE	10	9.53	1		95	71 - 130%
156-60-5	TRANS-1,2-DICHLOROETHENE	10	9.95	1		100	82 - 117%
1634-04-4	METHYL TERTIARY BUTYL ETHER	20	17.9	1		89	77 - 119%
75-34-3	1,1-DICHLOROETHANE	10	10.2	1		102	83 - 119%
108-05-4	VINYL ACETATE	10	9.43	2		94	76 - 121%
156-59-2	CIS-1,2-DICHLOROETHENE	10	9.93	1		99	83 - 117%
78-93-3	2-BUTANONE	40	36.2	10		91	70 - 135%
74-97-5	BROMOCHLOROMETHANE	10	10.1	1		101	83 - 121%
67-66-3	CHLOROFORM	10	9.5	1		95	82 - 119%
71-55-6	1,1,1-TRICHLOROETHANE	10	9.9	1		99	80 - 120%
594-20-7	2,2-DICHLOROPROPANE	10	10.3	1		103	83 - 125%
56-23-5	CARBON TETRACHLORIDE	10	9.91	1		99	77 - 122%
563-58-6	1,1-DICHLOROPROPENE	10	10.3	1		103	84 - 118%
107-06-2	1,2-DICHLOROETHANE	10	8.97	1		90	74 - 128%
71-43-2	BENZENE	10	9.97	1		100	83 - 117%

Data Package ID: VL1010128-1

Date Printed: Tuesday, October 26, 2010

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GC/MS Volatiles

Method SW8260_2514

Laboratory Control Sample and Laboratory Control Sample Duplicate

Lab Name: ALS Environmental -- FC

Work Order Number: 1010128

Client Name: Colorado Oil & Gas Conservation Commission

ClientProject ID: Complaint 200272771

Lab ID: VL101011-2LCS

Sample Matrix: WATER

% Moisture: N/A

Date Collected: N/A

Date Extracted: 10/11/2010

Date Analyzed: 10/11/2010

Prep Method: SW5030C

Prep Batch: VL101011-2

QCBatchID: VL101011-2-3

Run ID: VL101011-2A

Cleanup: NONE

Basis: N/A

File Name: B68261

Sample Aliquot: 10 ml

Final Volume: 10 ml

Result Units: UG/L

Clean DF: 1

CASNO	Target Analyte	Spike Added	LCS Result	Reporting Limit	Result Qualifier	LCS % Rec.	Control Limits
79-01-6	TRICHLOROETHENE	10	10	1		100	83 - 117%
78-87-5	1,2-DICHLOROPROPANE	10	10.4	1		104	84 - 120%
74-95-3	DIBROMOMETHANE	10	9.03	1		90	79 - 122%
75-27-4	BROMODICHLOROMETHANE	10	9.57	1		96	76 - 122%
10061-01-5	CIS-1,3-DICHLOROPROPENE	10	10.4	1		104	81 - 120%
108-10-1	4-METHYL-2-PENTANONE	40	36.7	10		92	73 - 125%
108-88-3	TOLUENE	10	9.75	1		98	82 - 113%
10061-02-6	TRANS-1,3-DICHLOROPROPENE	10	9.46	1		95	81 - 114%
79-00-5	1,1,2-TRICHLOROETHANE	10	9.05	1		91	78 - 116%
591-78-6	2-HEXANONE	40	32.7	10		82	71 - 124%
127-18-4	TETRACHLOROETHENE	10	10	1		100	81 - 115%
142-28-9	1,3-DICHLOROPROPANE	10	9.51	1		95	80 - 115%
124-48-1	DIBROMOCHLOROMETHANE	10	9.44	1		94	76 - 115%
106-93-4	1,2-DIBROMOETHANE	10	9.32	1		93	79 - 114%
544-10-5	1-CHLOROHEXANE	10	10.3	1		103	80 - 117%
108-90-7	CHLOROBENZENE	10	9.92	1		99	81 - 113%
630-20-6	1,1,1,2-TETRACHLOROETHANE	10	9.66	1		97	78 - 113%
100-41-4	ETHYLBENZENE	10	9.95	1		100	81 - 113%
136777-61-2	M+P-XYLENE	20	20.3	1		101	82 - 115%
95-47-6	O-XYLENE	10	10.5	1		105	81 - 115%
100-42-5	STYRENE	10	9.93	1		99	78 - 118%
75-25-2	BROMOFORM	10	8.56	1		86	70 - 120%
98-82-8	ISOPROPYLBENZENE	10	10.2	1		102	80 - 113%
96-18-4	1,2,3-TRICHLOROPROPANE	10	8.75	1		88	78 - 117%
79-34-5	1,1,2,2-TETRACHLOROETHANE	10	8.89	1		89	75 - 121%
108-86-1	BROMOBENZENE	10	9.09	1		91	81 - 114%
103-65-1	N-PROPYLBENZENE	10	10.2	1		102	79 - 116%

Data Package ID: VL1010128-1

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GC/MS Volatiles

Method SW8260_2514

Laboratory Control Sample and Laboratory Control Sample Duplicate

Lab Name: ALS Environmental -- FC

Work Order Number: 1010128

Client Name: Colorado Oil & Gas Conservation Commission

ClientProject ID: Complaint 200272771

Lab ID: VL101011-2LCS

Sample Matrix: WATER

% Moisture: N/A

Date Collected: N/A

Date Extracted: 10/11/2010

Date Analyzed: 10/11/2010

Prep Method: SW5030C

Prep Batch: VL101011-2

QCBatchID: VL101011-2-3

Run ID: VL101011-2A

Cleanup: NONE

Basis: N/A

File Name: B68261

Sample Aliquot: 10 ml

Final Volume: 10 ml

Result Units: UG/L

Clean DF: 1

CASNO	Target Analyte	Spike Added	LCS Result	Reporting Limit	Result Qualifier	LCS % Rec.	Control Limits
95-49-8	2-CHLOROTOLUENE	10	10.1	1		101	79 - 116%
108-67-8	1,3,5-TRIMETHYLBENZENE	10	10.1	1		101	78 - 116%
106-43-4	4-CHLOROTOLUENE	10	9.71	1		97	78 - 115%
98-06-6	TERT-BUTYLBENZENE	10	10.7	1		107	76 - 120%
95-63-6	1,2,4-TRIMETHYLBENZENE	10	9.73	1		97	80 - 117%
135-98-8	SEC-BUTYLBENZENE	10	10.4	1		104	78 - 115%
541-73-1	1,3-DICHLOROBENZENE	10	9.86	1		99	79 - 115%
99-87-6	P-ISOPROPYLTOLUENE	10	10.3	1		103	77 - 116%
106-46-7	1,4-DICHLOROBENZENE	10	9.44	1		94	82 - 114%
104-51-8	N-BUTYLBENZENE	10	10.7	1		107	79 - 117%
95-50-1	1,2-DICHLOROBENZENE	10	9.55	1		95	82 - 114%
96-12-8	1,2-DIBROMO-3-CHLOROPROPANE	10	8.42	2		84	73 - 125%
120-82-1	1,2,4-TRICHLOROBENZENE	10	10.3	1		103	75 - 120%
87-68-3	HEXACHLOROBUTADIENE	10	11.3	1		113	71 - 124%
91-20-3	NAPHTHALENE	10	9.63	1		96	71 - 131%
87-61-6	1,2,3-TRICHLOROBENZENE	10	10.4	1		104	70 - 131%

Data Package ID: VL1010128-1

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GC/MS Volatiles

Method SW8260_2514

Laboratory Control Sample and Laboratory Control Sample Duplicate

Lab Name: ALS Environmental -- FC

Work Order Number: 1010128

Client Name: Colorado Oil & Gas Conservation Commission

ClientProject ID: Complaint 200272771

Lab ID: VL101011-2LCSD

Sample Matrix: WATER

% Moisture: N/A

Date Collected: N/A

Date Extracted: 10/11/2010

Date Analyzed: 10/11/2010

Prep Method: SW5030C

Prep Batch: VL101011-2

QC Batch ID: VL101011-2-3

Run ID: VL101011-2A

Cleanup: NONE

Basis: N/A

File Name: B68262

Sample Aliquot: 10 ml

Final Volume: 10 ml

Result Units: UG/L

Clean DF: 1

CASNO	Target Analyte	Spike Added	LCSD Result	Reporting Limit	Result Qualifier	LCSD % Rec.	RPD Limit	RPD
75-71-8	DICHLORODIFLUOROMETHANE	10	9.18	1		92	20	4
74-87-3	CHLOROMETHANE	10	10.5	1		105	20	1
75-01-4	VINYL CHLORIDE	10	11.6	1		116	20	3
74-83-9	BROMOMETHANE	10	10.4	1		104	20	0
75-00-3	CHLOROETHANE	10	9.51	1		95	20	0
75-69-4	TRICHLOROFLUOROMETHANE	10	9.39	1		94	20	4
75-35-4	1,1-DICHLOROETHENE	10	9.38	1		94	20	2
76-13-1	1,1,2-TRICHLORO-1,2,2-TRIFLUOROETHA	10	10	1		100	20	3
67-64-1	ACETONE	40	33	10		83	30	5
74-88-4	IODOMETHANE	10	9.58	1		96	20	4
75-15-0	CARBON DISULFIDE	10	10	1		100	20	4
75-09-2	METHYLENE CHLORIDE	10	9.13	1		91	20	4
156-60-5	TRANS-1,2-DICHLOROETHENE	10	9.78	1		98	20	2
1634-04-4	METHYL TERTIARY BUTYL ETHER	20	18	1		90	20	1
75-34-3	1,1-DICHLOROETHANE	10	10.2	1		102	20	0
108-05-4	VINYL ACETATE	10	10.1	2		101	20	7
156-59-2	CIS-1,2-DICHLOROETHENE	10	9.6	1		96	20	3
78-93-3	2-BUTANONE	40	37.1	10		93	30	3
74-97-5	BROMOCHLOROMETHANE	10	9.61	1		96	20	5
67-66-3	CHLOROFORM	10	9.46	1		95	20	0
71-55-6	1,1,1-TRICHLOROETHANE	10	9.52	1		95	20	4
594-20-7	2,2-DICHLOROPROPANE	10	9.78	1		98	20	6
56-23-5	CARBON TETRACHLORIDE	10	9.41	1		94	20	5
563-58-6	1,1-DICHLOROPROPENE	10	9.97	1		100	20	3
107-06-2	1,2-DICHLOROETHANE	10	8.87	1		89	20	1
71-43-2	BENZENE	10	10	1		100	20	0
79-01-6	TRICHLOROETHENE	10	9.79	1		98	20	3

Data Package ID: VL1010128-1

Date Printed: Tuesday, October 26, 2010

ALS Environmental -- FC

Page 4 of 6

LIMS Version: 6.423A

GC/MS Volatiles

Method SW8260_2514

Laboratory Control Sample and Laboratory Control Sample Duplicate

Lab Name: ALS Environmental -- FC

Work Order Number: 1010128

Client Name: Colorado Oil & Gas Conservation Commission

ClientProject ID: Complaint 200272771

Lab ID: VL101011-2LCSD

Sample Matrix: WATER

% Moisture: N/A

Date Collected: N/A

Date Extracted: 10/11/2010

Date Analyzed: 10/11/2010

Prep Method: SW5030C

Prep Batch: VL101011-2

QCBatchID: VL101011-2-3

Run ID: VL101011-2A

Cleanup: NONE

Basis: N/A

File Name: B68262

Sample Aliquot: 10 ml

Final Volume: 10 ml

Result Units: UG/L

Clean DF: 1

CASNO	Target Analyte	Spike Added	LCSD Result	Reporting Limit	Result Qualifier	LCSD % Rec.	RPD Limit	RPD
78-87-5	1,2-DICHLOROPROPANE	10	10.4	1		104	20	1
74-95-3	DIBROMOMETHANE	10	9.13	1		91	20	1
75-27-4	BROMODICHLOROMETHANE	10	9.69	1		97	20	1
10061-01-5	CIS-1,3-DICHLOROPROPENE	10	10.4	1		104	20	0
108-10-1	4-METHYL-2-PENTANONE	40	38.2	10		96	30	4
108-88-3	TOLUENE	10	10.2	1		102	20	4
10061-02-6	TRANS-1,3-DICHLOROPROPENE	10	9.81	1		98	20	4
79-00-5	1,1,2-TRICHLOROETHANE	10	9.54	1		95	20	5
591-78-6	2-HEXANONE	40	35.5	10		89	30	8
127-18-4	TETRACHLOROETHENE	10	10.4	1		104	20	3
142-28-9	1,3-DICHLOROPROPANE	10	9.92	1		99	20	4
124-48-1	DIBROMOCHLOROMETHANE	10	9.59	1		96	20	2
106-93-4	1,2-DIBROMOETHANE	10	9.66	1		97	20	4
544-10-5	1-CHLOROHEXANE	10	10.7	1		107	20	4
108-90-7	CHLOROBENZENE	10	10.2	1		102	20	3
630-20-6	1,1,1,2-TETRACHLOROETHANE	10	9.68	1		97	20	0
100-41-4	ETHYLBENZENE	10	9.95	1		99	20	0
136777-61-2	M+P-XYLENE	20	20.8	1		104	20	3
95-47-6	O-XYLENE	10	10.4	1		104	20	1
100-42-5	STYRENE	10	10.2	1		102	20	2
75-25-2	BROMOFORM	10	8.67	1		87	20	1
98-82-8	ISOPROPYLBENZENE	10	10.2	1		102	20	0
96-18-4	1,2,3-TRICHLOROPROPANE	10	9.13	1		91	20	4
79-34-5	1,1,2,2-TETRACHLOROETHANE	10	9.03	1		90	20	2
108-86-1	BROMOBENZENE	10	9.27	1		93	20	2
103-65-1	N-PROPYLBENZENE	10	10.2	1		102	20	0
95-49-8	2-CHLOROTOLUENE	10	10.3	1		103	20	2

Data Package ID: VL1010128-1

Date Printed: Tuesday, October 26, 2010

ALS Environmental -- FC

Page 5 of 6

LIMS Version: 6.423A

GC/MS Volatiles

Method SW8260_2514

Laboratory Control Sample and Laboratory Control Sample Duplicate

Lab Name: ALS Environmental -- FC

Work Order Number: 1010128

Client Name: Colorado Oil & Gas Conservation Commission

ClientProject ID: Complaint 200272771

Lab ID: VL101011-2LCSD

Sample Matrix: WATER

% Moisture: N/A

Date Collected: N/A

Date Extracted: 10/11/2010

Date Analyzed: 10/11/2010

Prep Method: SW5030C

Prep Batch: VL101011-2

QC Batch ID: VL101011-2-3

Run ID: VL101011-2A

Cleanup: NONE

Basis: N/A

File Name: B68262

Sample Aliquot: 10 ml

Final Volume: 10 ml

Result Units: UG/L

Clean DF: 1

CASNO	Target Analyte	Spike Added	LCSD Result	Reporting Limit	Result Qualifier	LCSD % Rec.	RPD Limit	RPD
108-67-8	1,3,5-TRIMETHYLBENZENE	10	10	1		100	20	0
106-43-4	4-CHLOROTOLUENE	10	10.1	1		101	20	4
98-06-6	TERT-BUTYLBENZENE	10	10.7	1		107	20	0
95-63-6	1,2,4-TRIMETHYLBENZENE	10	9.74	1		97	20	0
135-98-8	SEC-BUTYLBENZENE	10	10.5	1		105	20	1
541-73-1	1,3-DICHLOROBENZENE	10	9.87	1		99	20	0
99-87-6	P-ISOPROPYLTOLUENE	10	10.2	1		102	20	1
106-46-7	1,4-DICHLOROBENZENE	10	9.41	1		94	20	0
104-51-8	N-BUTYLBENZENE	10	10.6	1		106	20	1
95-50-1	1,2-DICHLOROBENZENE	10	9.36	1		94	20	2
96-12-8	1,2-DIBROMO-3-CHLOROPROPANE	10	8.51	2		85	20	1
120-82-1	1,2,4-TRICHLOROBENZENE	10	10.4	1		104	20	2
87-68-3	HEXACHLOROBUTADIENE	10	11	1		110	20	2
91-20-3	NAPHTHALENE	10	9.99	1		100	20	4
87-61-6	1,2,3-TRICHLOROBENZENE	10	11.1	1		111	20	6

Surrogate Recovery LCS/LCSD

CASNO	Target Analyte	Spike Added	LCS % Rec.	LCS Flag	LCSD % Rec.	LCSD Flag	Control Limits
460-00-4	4-BROMOFLUOROBENZENE	25	97		98		87 - 108
1868-53-7	DIBROMOFLUOROMETHANE	25	100		96		89 - 112
2037-26-5	TOLUENE-D8	25	100		101		87 - 109

Data Package ID: VL1010128-1

Date Printed: Tuesday, October 26, 2010

ALS Environmental -- FC

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LIMS Version: 6.423A

Data File : C:\HPCHEM\1\DATA\101110\B68264.D

Vial: 5

Acq On : 11 Oct 2010 9:00

Operator: twk-sop525r14

Sample : VL101011-2MB

Inst : CSS Instr

Misc : 10mL un-heated purge (water)

Multiplr: 1.00

MS Integration Params: ettics.p

Quant Time: Oct 11 10:38 2010

Quant Results File: 092410W.RES

Quant Method : C:\HPCHEM\1\METHODS\092410W.M (RTE Integrator)

Title : GC/MS Volatiles (S.O.P. 525)

Last Update : Mon Oct 11 10:37:17 2010

Response via : Initial Calibration

DataAcq Meth : 092410W

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Fluorobenzene	9.49	96	2154478	25.00	ppb	0.00
56) Chlorobenzene-d5	12.50	82	808053	25.00	ppb	0.00
78) 1,4-Dichlorobenzene-d4	14.43	152	533986	25.00	ppb	0.00

System Monitoring Compounds

37) Dibromofluoromethane	8.74	113	604694	23.78	ppb	0.00
Spiked Amount	25.000	Range	89 - 112	Recovery	=	95.12%
41) 1,2-dichloroethane-d4	9.19	65	428971	22.72	ppb	0.00
Spiked Amount	25.000	Range	80 - 125	Recovery	=	90.88%
57) Toluene-d8	11.11	98	1918535	25.50	ppb	0.00
Spiked Amount	25.000	Range	87 - 109	Recovery	=	102.00%
77) 4-Bromofluorobenzene	13.49	176	412274	23.32	ppb	0.00
Spiked Amount	25.000	Range	87 - 108	Recovery	=	93.28%

Target Compounds

						Qvalue
13) Acetone	6.03	43	4135	Below Cal		87
16) Allyl chloride	6.29	76	3310	0.19 ppb	# No	1
18) Methylene chloride	6.61	84	3466	Below Cal	#	61

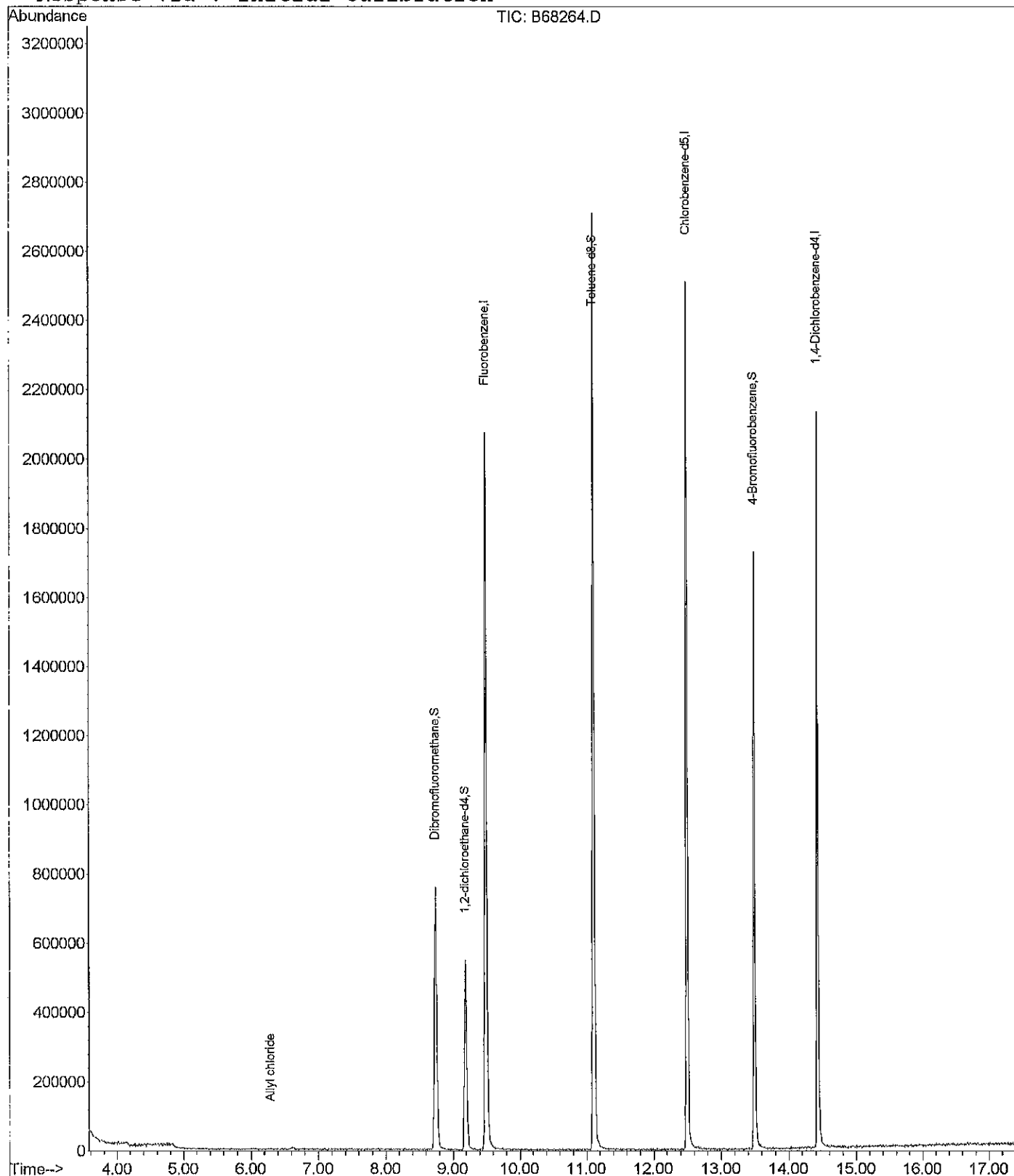
Quantitation Report

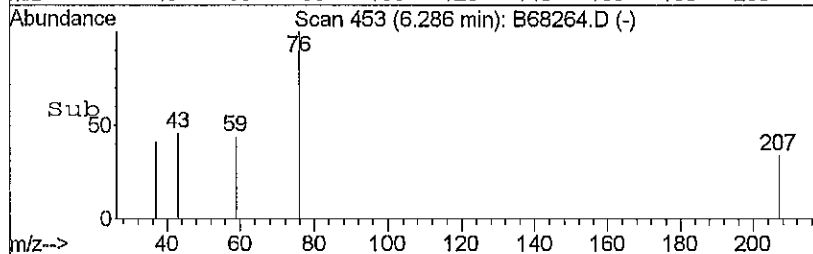
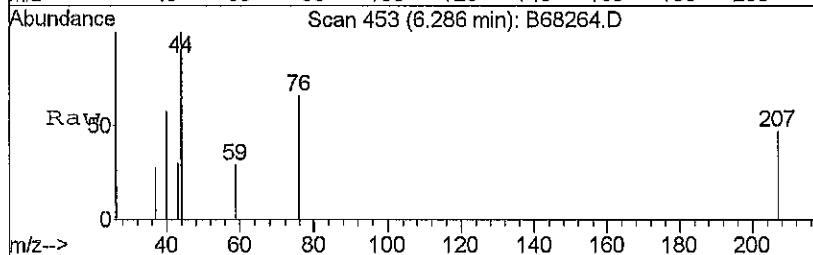
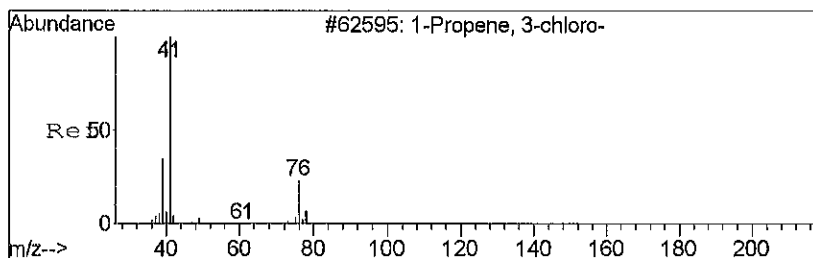
Data File : C:\HPCHEM\1\DATA\101110\B68264.D
Acq On : 11 Oct 2010 9:00
Sample : VL101011-2MB
Misc : 10mL un-heated purge (water)
MS Integration Params: ettics.p
Quant Time: Oct 11 10:38 2010

Vial: 5
Operator: twk-sop525r14
Inst : CSS Instr
Multiplr: 1.00

Quant Results File: 092410W.RES

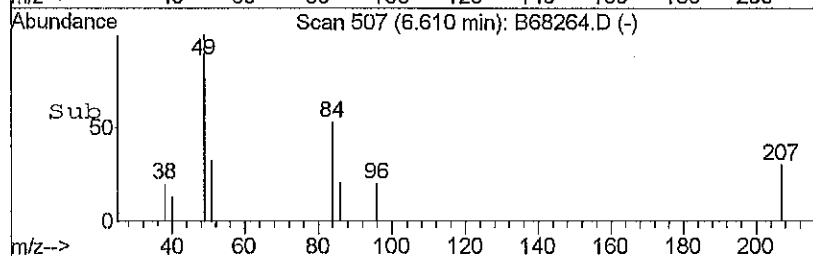
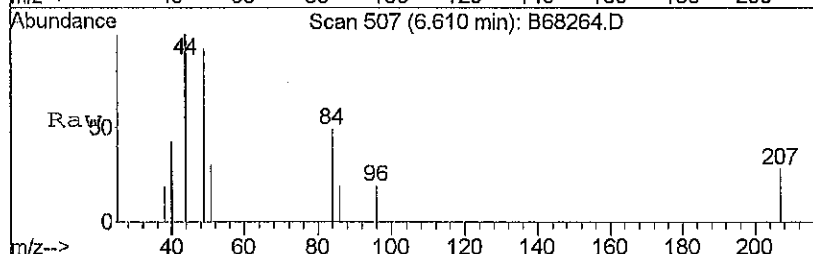
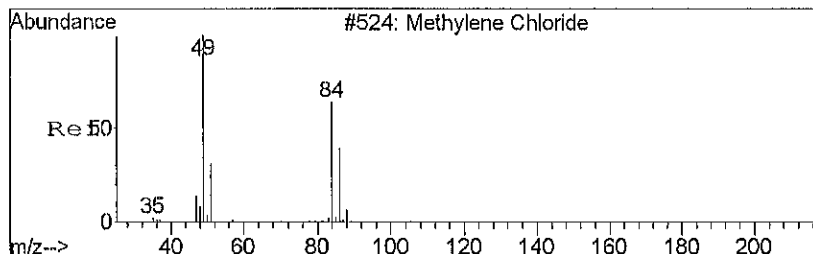
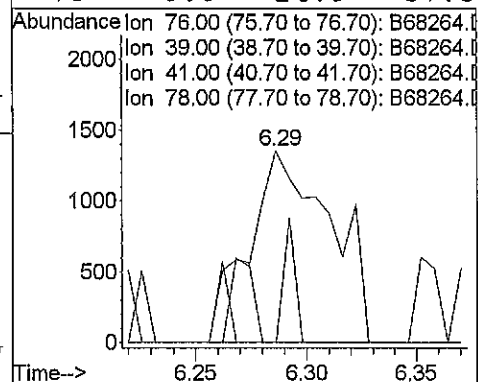
Method : C:\HPCHEM\1\METHODS\092410W.M (RTE Integrator)
Title : GC/MS Volatiles (S.O.P. 525)
Last Update : Mon Oct 11 10:37:17 2010
Response via : Initial Calibration





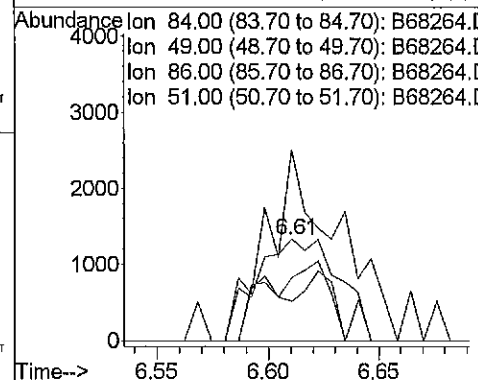
#16
 Allyl chloride
 Concen: 0.19 ppb
 RT: 6.29 min Scan# 453
 Delta R.T. -0.12 min
 Lab File: B68264.D
 Acq: 11 Oct 2010 9:00

Tgt Ion: 76 Resp: 3310
 Ion Ratio Lower Upper
 76 100
 39 0.0 118.1 177.1#
 41 0.0 200.3 300.5#
 78 0.0 24.9 37.3#



#18
 Methylene chloride
 Concen: Below Cal
 RT: 6.61 min Scan# 507
 Delta R.T. -0.00 min
 Lab File: B68264.D
 Acq: 11 Oct 2010 9:00

Tgt Ion: 84 Resp: 3466
 Ion Ratio Lower Upper
 84 100
 49 188.7 81.2 189.4
 86 39.3 37.5 87.5
 51 62.2 24.5 57.1#



Data File : C:\HPCHEM\1\DATA\101110\B68264.D

Vial: 5

Acq On : 11 Oct 2010 9:00

Operator: twk-sop525r14

Sample : VL101011-2MB

Inst : CSS Instr

Misc : 10mL un-heated purge (water)

Multiplr: 1.00

MS Integration Params: ettics.p

Quant Time: Oct 13 16:38 2010

Quant Results File: 110309WH.RES

Quant Method : C:\HPCHEM\1\METHODS\110309WH.M (RTE Integrator)

Title : GC/MS Volatiles (S.O.P. 525)

Last Update : Tue Aug 10 13:51:21 2010

Response via : Initial Calibration

DataAcq Meth : 092410W

Internal Standards	R.T.	Q	Ion	Response	Conc	Units	Dev (Min)
1) Fluorobenzene	9.49	96		2154478	25.00	ppb	-0.02

Target Compounds

Qvalue

m 10/13/10

Data File : C:\HPCHEM\1\DATA\101110\B68278.D

Vial: 17

Acq On : 11 Oct 2010 14:14

Operator: twk-sop525r14

Sample : 1010128-1

Inst : CSS Instr

Misc : 10mL un-heated purge (water)

Multiplr: 1.00

MS Integration Params: ettics.p

Quant Time: Oct 11 14:46 2010

Quant Results File: 092410W.RES

Quant Method : C:\HPCHEM\1\METHODS\092410W.M (RTE Integrator)

Title : GC/MS Volatiles (S.O.P. 525)

Last Update : Mon Oct 11 10:37:17 2010

Response via : Initial Calibration

DataAcq Meth : 092410W

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Fluorobenzene	9.50	96	1978970	25.00	ppb	0.00
56) Chlorobenzene-d5	12.50	82	785378	25.00	ppb	0.00
78) 1,4-Dichlorobenzene-d4	14.44	152	515433	25.00	ppb	0.00

System Monitoring Compounds

37) Dibromofluoromethane	8.75	113	579654	24.82	ppb	0.00
Spiked Amount	25.000	Range	89 - 112	Recovery	=	99.28%
41) 1,2-dichloroethane-d4	9.19	65	403446	23.26	ppb	0.00
Spiked Amount	25.000	Range	80 - 125	Recovery	=	93.04%
57) Toluene-d8	11.11	98	1798159	24.59	ppb	0.00
Spiked Amount	25.000	Range	87 - 109	Recovery	=	98.36%
77) 4-Bromofluorobenzene	13.50	176	411484	23.95	ppb	0.00
Spiked Amount	25.000	Range	87 - 108	Recovery	=	95.80%

Target Compounds

Qvalue

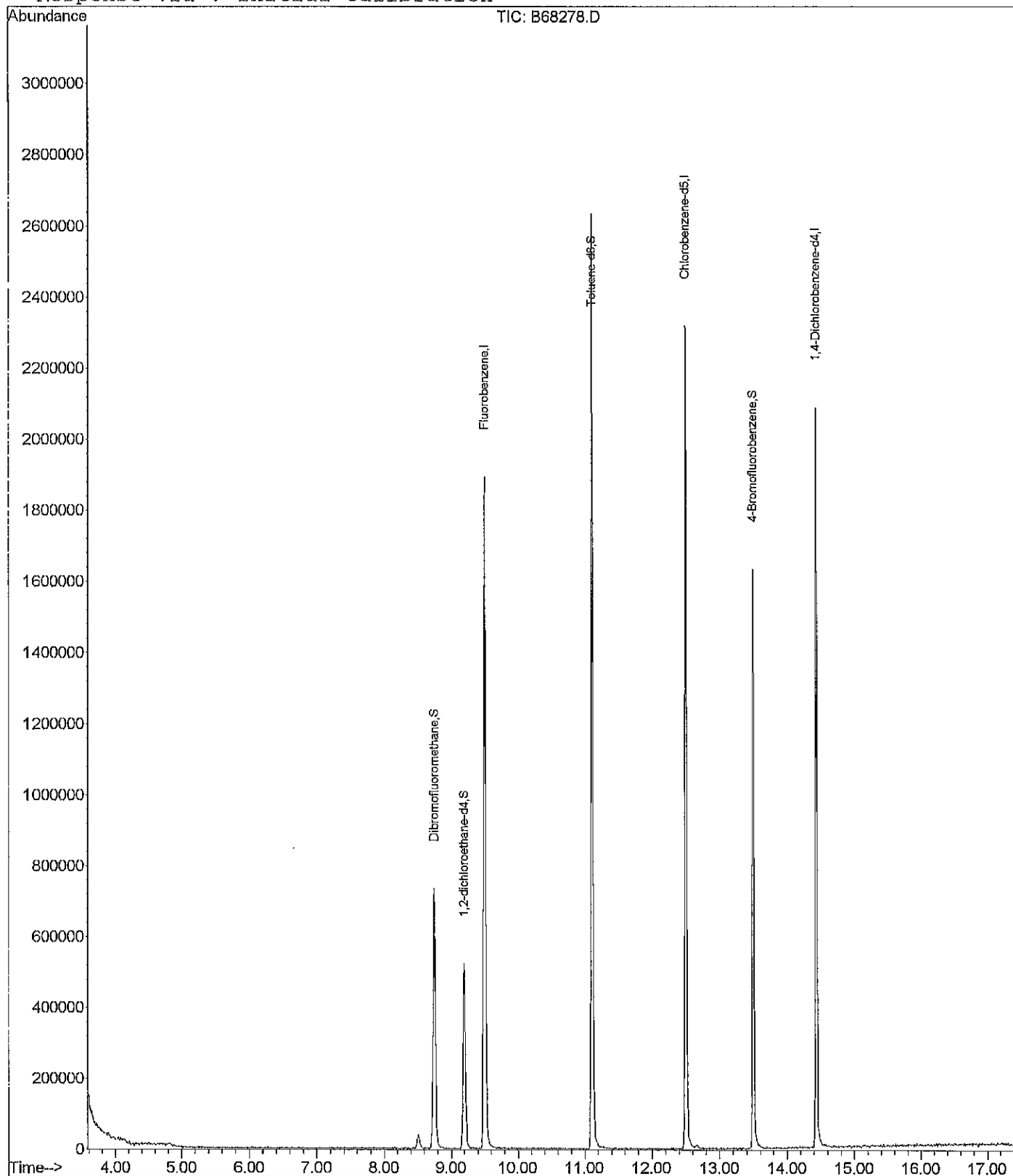
Quantitation Report

Data File : C:\HPCHEM\1\DATA\101110\B68278.D
 Acq On : 11 Oct 2010 14:14
 Sample : 1010128-1
 Misc : 10mL un-heated purge (water)
 MS Integration Params: ettics.p
 Quant Time: Oct 11 14:46 2010

Vial: 17
 Operator: twk-sop525r14
 Inst : CSS Instr
 Multiplr: 1.00

Quant Results File: 092410W.RES

Method : C:\HPCHEM\1\METHODS\092410W.M (RTE Integrator)
 Title : GC/MS Volatiles (S.O.P. 525)
 Last Update : Mon Oct 11 10:37:17 2010
 Response via : Initial Calibration



Data File : C:\HPCHEM\1\DATA\101110\B68278.D

Acq On : 11 Oct 2010 14:14

Sample : 1010128-1

Misc : 10mL un-heated purge (water)

MS Integration Params: ettics.p

Vial: 17

Operator: twk-sop525r14

Inst : CSS Instr

Multiplr: 1.00

Quant Method : C:\HPCHEM\1\METHODS\092410W.M (RTE Integrator)

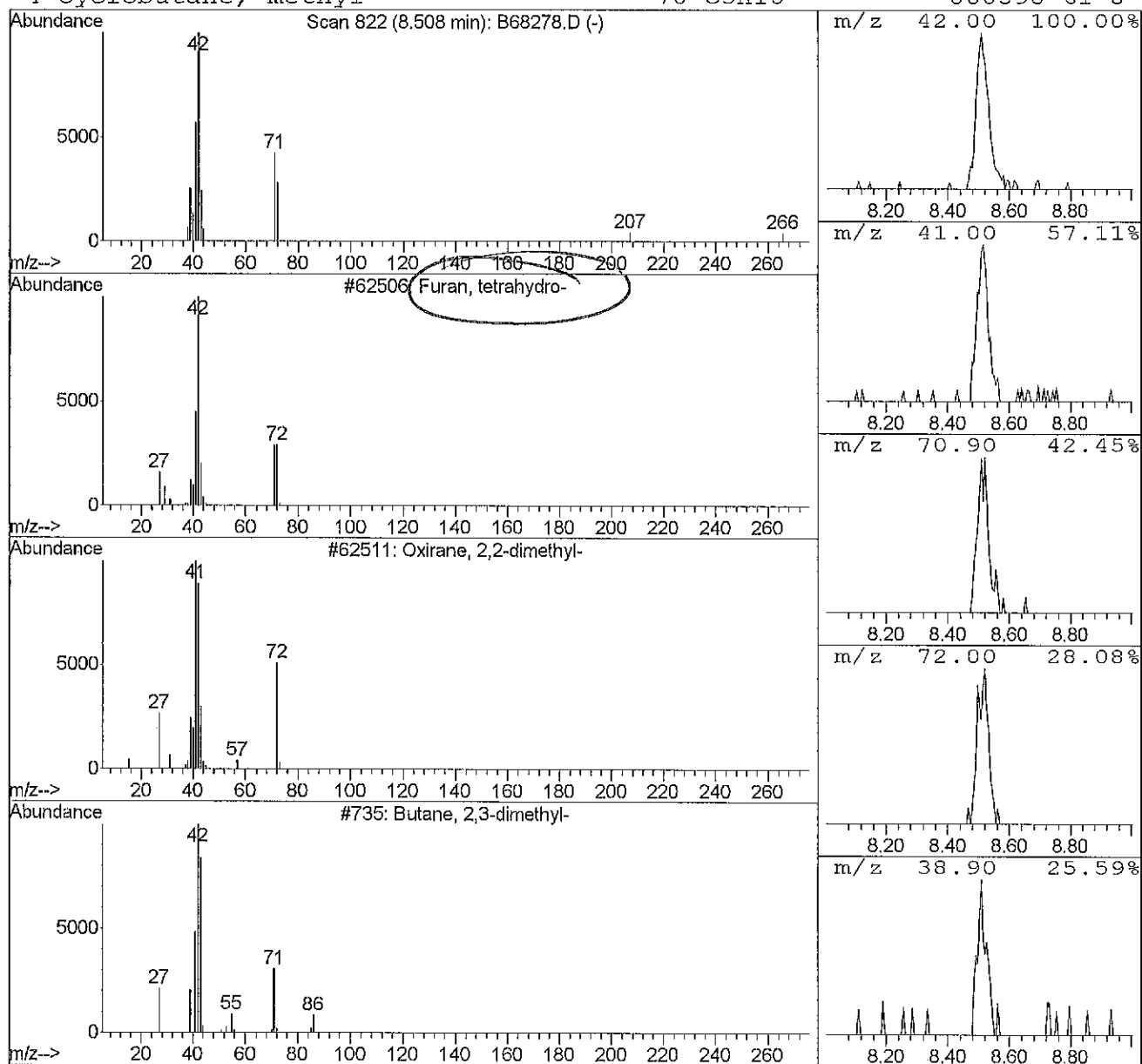
Title : GC/MS Volatiles (S.O.P. 525)

Library : C:\DATABASE\NBS75K.L

Peak Number 1 Furan, tetrahydro- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.51	0.67 ppb	111086	Fluorobenzene	9.50

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Furan, tetrahydro-	72	C4H8O	000109-99-9	78
2			Oxirane, 2,2-dimethyl-	72	C4H8O	000558-30-5	9
3			Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	9
4			Cyclobutane, methyl-	70	C5H10	000598-61-8	9



Data File : C:\HPCHEM\1\DATA\101110\B68279.D

Vial: 17

Acq On : 11 Oct 2010 14:36

Operator: twk-sop525r14

Sample : 1010128-2

Inst : CSS Instr

Misc : 10mL un-heated purge (water)

Multiplr: 1.00

MS Integration Params: ettics.p

Quant Time: Oct 11 15:02 2010

Quant Results File: 092410W.RES

Quant Method : C:\HPCHEM\1\METHODS\092410W.M (RTE Integrator)

Title : GC/MS Volatiles (S.O.P. 525)

Last Update : Mon Oct 11 10:37:17 2010

Response via : Initial Calibration

DataAcq Meth : 092410W

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Fluorobenzene	9.50	96	1816382	25.00	ppb	0.00
56) Chlorobenzene-d5	12.50	82	704928	25.00	ppb	0.00
78) 1,4-Dichlorobenzene-d4	14.43	152	468075	25.00	ppb	0.00

System Monitoring Compounds

37) Dibromofluoromethane	8.75	113	501558	23.39	ppb	0.00
Spiked Amount 25.000	Range 89 - 112		Recovery =	93.56%		
41) 1,2-dichloroethane-d4	9.19	65	377271	23.70	ppb	0.00
Spiked Amount 25.000	Range 80 - 125		Recovery =	94.80%		
57) Toluene-d8	11.11	98	1639458	24.98	ppb	0.00
Spiked Amount 25.000	Range 87 - 109		Recovery =	99.92%		
77) 4-Bromofluorobenzene	13.49	176	369569	23.97	ppb	0.00
Spiked Amount 25.000	Range 87 - 108		Recovery =	95.88%		

Target Compounds

18) Methylene chloride	6.62	84	21395	Below Cal	Qvalue 90
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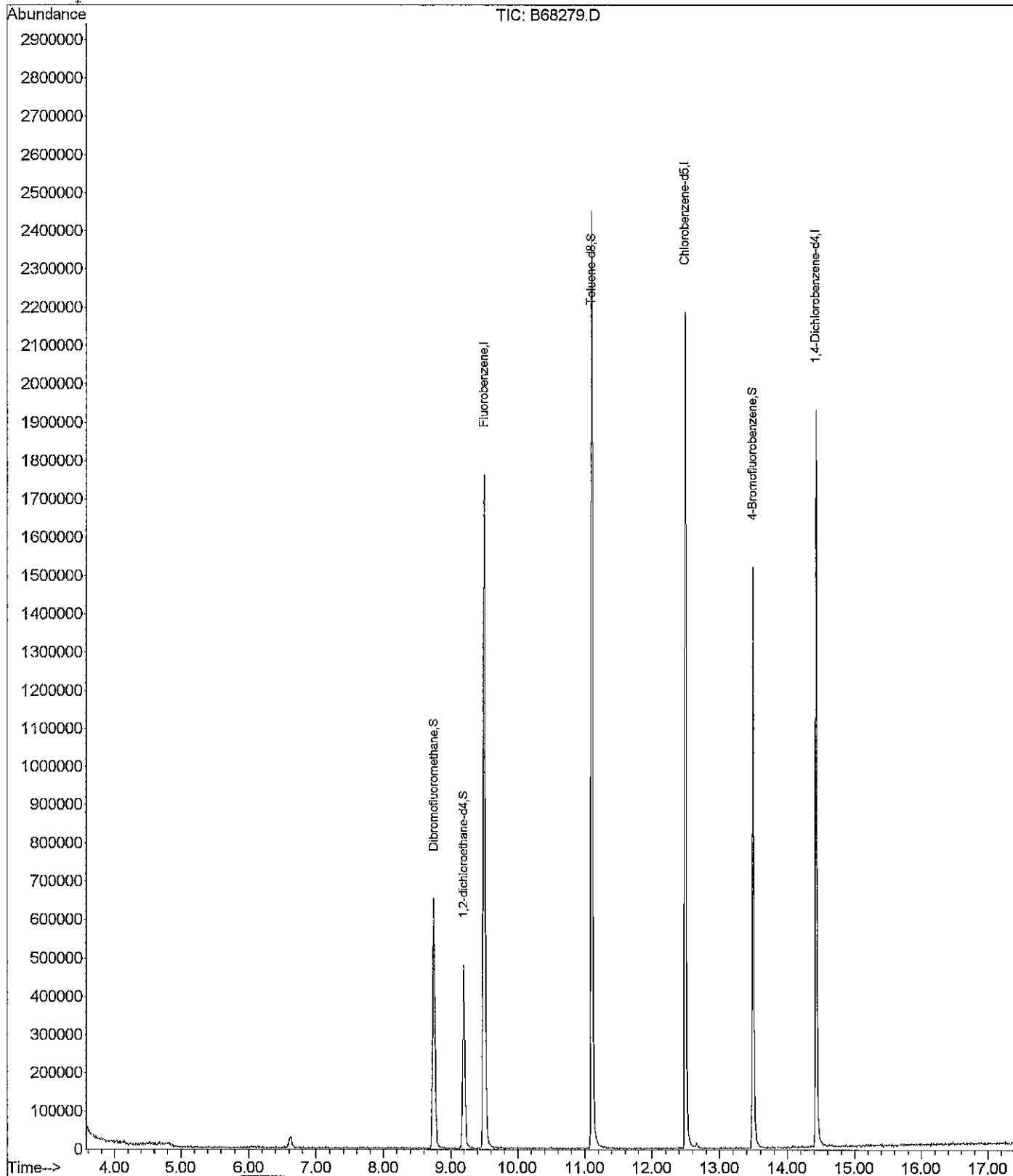
Quantitation Report

Data File : C:\HPCHEM\1\DATA\101110\B68279.D
 Acq On : 11 Oct 2010 14:36
 Sample : 1010128-2
 Misc : 10mL un-heated purge (water)
 MS Integration Params: ettics.p
 Quant Time: Oct 11 15:02 2010

Vial: 17
 Operator: twk-sop525r14
 Inst : CSS Instr
 Multiplr: 1.00

Quant Results File: 092410W.RES

Method : C:\HPCHEM\1\METHODS\092410W.M (RTE Integrator)
 Title : GC/MS Volatiles (S.O.P. 525)
 Last Update : Mon Oct 11 10:37:17 2010
 Response via : Initial Calibration



Tentatively Identified Compound (LSC) summary

Operator ID: twk-sop525r14 Date Acquired: 11 Oct 2010 14:36
 Data File: C:\HPCHEM\1\DATA\101110\B68279.D
 Name: 1010128-2
 Misc: 10mL un-heated purge (water)
 Method: C:\HPCHEM\1\METHODS\092410W.M (RTE Integrator)
 Title: GC/MS Volatiles (S.O.P. 525)
 Library Searched: C:\DATABASE\NBS75K.L

TIC Top Hit name	RT	EstConc	Units	Area	IntStd	ISRT	ISArea	ISConc

B68279.D 092410W.M		Wed Oct 13	16:39:16	2010				