



SHIPPING DOCUMENTS

Client Contact Information				Bottle Order ID : B2504002				Courier : FGaldun				1 of 2 COCs																	
Client ID : GFEL01 Project ID : to Emerald St.				Sampler Name(s) : FRANK Galdun				Analysis				Matrix																	
Customer Name : GFE LLC Address : 58 Nokomis Ave City : Lake Hiawatha State : NJ Zip Code : 07034 Country :				Project Manager : Frank galdun				AIR ANALYSIS CHAIN-OF-CUSTODY Batch Certified																					
				Phone Number : 646-542-3465																									
				Fax Number : 973-334-1692																									
				Site Details: 416 CLINTON ST. HEMPSTEAD, NY																									
				Analysis Turnaround Time																									
				Standard : 10 business days OR				Data Package Type SAME PACKAGE AS PREVIOUS ROUND																					
				Rush (Specify): Days				EDD Type :																					
Sample Identification	Sample Date(s)	Time Start (24 hr Clock)	Time Stop (24 hr Clock)	Can Vacuum in Field ("Hg) (Start)	Can Vacuum in Field ("Hg) (Stop)**	Interior Temp. (F) (Start)	Interior Temp. (F) (Stop)	Out going Can Pressure ("Hg)(Lab)	In coming Can Pressure ("Hg)(Lab)	Flow Reg. ID	Can ID		Flow Controller Readout (ml/min)	Can Cert ID	TO-15	Indoor/Ambient Air	Soil Gas												
SS2R	4/12/25	7:00	7:00	0.0	0.0	50	52	-30	-2.2	10648	10311	6 L	50	VL042217.D	(		(												
Temperature (Fahrenheit) <table border="1" style="width:100%; border-collapse: collapse;"> <tr> <td></td> <td>Ambient</td> <td>Maximum</td> <td>Minimum</td> </tr> <tr> <td>Start</td> <td></td> <td></td> <td></td> </tr> <tr> <td>Stop</td> <td></td> <td></td> <td></td> </tr> </table> GC/MS Analyst Signature (TO-15)											Ambient	Maximum	Minimum	Start				Stop				** Submittal of this COC indicates approval of the analysis based on existing conditions. Please follow the instructions on the back of this COC.							
	Ambient	Maximum	Minimum																										
Start																													
Stop																													
Pressure (Inches of Hg) <table border="1" style="width:100%; border-collapse: collapse;"> <tr> <td></td> <td>Ambient</td> <td>Maximum</td> <td>Minimum</td> </tr> <tr> <td>Start</td> <td></td> <td></td> <td></td> </tr> <tr> <td>Stop</td> <td></td> <td></td> <td></td> </tr> </table>											Ambient	Maximum	Minimum	Start				Stop											
	Ambient	Maximum	Minimum																										
Start																													
Stop																													
Special Instructions/QC Requirements & Comments :																													
Suspected Contamination: High Medium <u>Low</u> PID Readings: <u>0.0</u> Sampling site (State):																													
Quick Connector required : <u>NO</u>																													
Canisters Shipped by: <u>SEM</u>				Date/Time: <u>04/07/25</u>				Canisters Received by: <u>AG</u>				Date/Time: <u>4/14/25 0700</u>																	
Samples Relinquished by: <u>SEM</u>				Date/Time: <u>4/12/25</u>				Received by:				Date/Time:																	
Relinquished by:				Date/Time:				Received by:				Date/Time:																	

Client Contact Information					Bottle Order ID : B2504002					Courier : FGALDUN					2 of 2 COCs				
Client ID : GFEL01 Project ID : 10 Eldridge St.					Sampler Name(s) : FRANK GALDUN					Analysis					Matrix				
Customer : GFE LLC Name : Address : 58 Nokomis Ave City : Lake Hiawatha State : NJ Zip Code : 07034 Country :					Project Manager : Frank galdun					AIR ANALYSIS CHAIN-OF-CUSTODY Batch Certified									
					Phone Number : 646-542-3465														
					Fax Number : 973-334-1692														
					Site Details: 416 CLINTON ST. HEMPSTEAD, NY														
					Analysis Turnaround Time					Data Package Type : SAME PACKAGE AS PREVIOUS ROUND					Indoor/Ambient Air Soil Gas				
					Standard : 10 business days OR					EDD Type :									
Country :					Rush (Specify): Days														
Sample Identification	Sample Date(s)	Time Start (24 hr Clock)	Time Stop (24 hr Clock)	Can Vacuum in Field ("Hg) (Start)	Can Vacuum in Field ("Hg) (Stop)**	Interior Temp. (F) (Start)	Interior Temp. (F) (Stop)	Out going Can Pressure ("Hg)(Lab)	In coming Can Pressure ("Hg)(Lab)	Flow Reg. ID	Can ID		Flow Controller Readout (ml/min)	Can Cert ID	TO-15				
553	4/12/25	7:02	8:02	OVER 30	1	65	67	-30	-3.1	10226	10606	6 L	50	VL042217.D	1				
Temperature (Fahrenheit)										GC/MS Analyst Signature (TO-15)									
		Ambient		Maximum		Minimum													
Start																			
Stop																			
Pressure (Inches of Hg)										** Submittal of this COC indicates approval of the analysis based on existing conditions. Please follow the instructions on the back of this COC.									
		Ambient		Maximum		Minimum													
Start																			
Stop																			
Special Instructions/QC Requirements & Comments :																			
Suspected Contamination: High Medium <u>Low</u> PID Readings: <u>0.0</u>																			
Sampling site (State):																			
Quick Connector required : <u>NO</u>																			
Canisters Shipped by: <u>SG</u>					Date/Time: <u>04/16/25</u>					Canisters Received by: <u>AS</u>					Date/Time: <u>4/14/25 0700</u>				
Samples Relinquished by: <u>AS</u>					Date/Time: <u>4/12/25</u>					Received by:					Date/Time:				
Relinquished by:					Date/Time:					Received by:					Date/Time:				
B2504002 - 6																			

Laboratory Certification

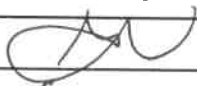
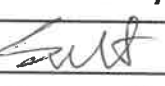
Certified By	License No.
CAS EPA CLP Contract	68HERH20D0011
Connecticut	PH-0830
DOD ELAP (ANAB)	L2219
Maine	2024021
Maryland	296
New Hampshire	255424 Rev 1
New Jersey	20012
New York	11376
Pennsylvania	68-00548
Soil Permit	525-24-234-08441
Texas	T104704488

Internal Chain of Custody

Instructions: Use 1 form for each 20 samples of aliquot

Laboratory Person Breaking Field Seal on Sample Shuttle & Accepting Responsibility for Sample			
Laboratory: <u>Chemtech</u>		Location: <u>284 Sheffield Street, Mountainside, NJ 7092</u>	
NOISE		Title: <u>Sample Custodian</u>	
Field Sample Seal No.: <u>Q1801</u>		Date Broken <u>4/14/2025</u> Military Time Seal Broken: <u>07:00:00</u>	
Case No.: <u>416 Clinton St Hempstead</u>		Analytical Parameter/Fraction <u>TO-15</u>	

Sample No.	Aliquot/Extract No.	Sample No.	Aliquot/Extract No.
Q1801-01	SS2R		
Q1801-02	SS3		

Date	Time	Relinquished By	Received By	Purpose of Change of Custody
4/14	1215	Signature 	Signature 	
		Printed Name <u>BOUSE N.</u>	Printed Name <u>Samuel J. J. J.</u>	
		Signature	Signature	
		Printed Name	Printed Name	
		Signature	Signature	
		Printed Name	Printed Name	
		Signature	Signature	
		Printed Name	Printed Name	
		Signature	Signature	
		Printed Name	Printed Name	
		Signature	Signature	
		Printed Name	Printed Name	
		Signature	Signature	
		Printed Name	Printed Name	
		Signature	Signature	
		Printed Name	Printed Name	

Distribution: White - Original (Sent With Report) Yellow - Contractor Archive Pink - Sample Custodian - Interim Copy

DATA REPORTING QUALIFIERS- ORGANIC

For reporting results, the following “ Results Qualifiers” are used:

Value	If the result is a value greater than or equal to the detection limit, report the value
U	Indicates the compound was analyzed for but was not detected. Report the minimum detection limit for the sample with the U, i.e. “10 U”. This is not necessarily the instrument detection limit attainable for this particular sample based on any concentration or dilution that may have been required.
ND	Indicates the analyte was analyzed for, but not detected
J	Indicates an estimated value. This flag is used: (1) When estimating a concentration for a tentatively identified compound (library search hits, where a 1:1 response is assumed.) (2) When the mass spectral data indicated the identification, however the result was less than the specified detection limit greater than zero. If the detection limit was 10ug/L and a concentration of 3 ug/L was calculated report as 3 J. This flag is used when similar situation arise on any organic parameter i.e. Pest, PCB and others.
B	Indicates the analyte was found in the blank as well as the sample report as “12 B”.
E	Indicates the analyte ‘s concentration exceeds the calibrated range of the instrument for that specific analysis.
D	This flag identifies all compounds identified in an analysis at a secondary dilution factor.
P	This flag is used for Pesticide/PCB target analyte when there is >25% difference for detected concentrations between the two GC columns. The lower of the two values is reported on Form 1 and flagged with a “P”.
N	This flag indicates presumptive evidence of a compound. This is only used for tentatively identified compounds (TICs), where the identification is based on a mass spectral library search. It applies to all TIC results. For generic characterization of a TIC, such as chlorinated hydrocarbon, the flag is not used.
A	This flag indicates that a Tentatively Identified Compound is a suspected aldol-condensation product.
Q	Indicates the LCS did not meet the control limits requirements

LAB CHRONICLE

OrderID: Q1801
Client: GFE LLC
Contact: Frank Galdun

OrderDate: 4/14/2025 11:40:00 AM
Project: 416 Clinton St Hempstead
Location: L31

LabID	ClientID	Matrix	Test	Method	Sample Date	Prep Date	Anal Date	Received
Q1801-01	SS2R	Air	TO-15	TO-15	04/12/25		04/21/25	04/14/25
Q1801-02	SS3	Air	TO-15	TO-15	04/12/25		04/21/25	04/14/25

Hit Summary Sheet
SW-846

SDG No.: Q1801
Client: GFE LLC

Sample ID	Client ID	Matrix	Parameter	Concentration	C	MDL	RDL	Units
Client ID:	SS2R							
Q1801-01	SS2R	Air	Heptane	1.89	J	1.80	8.20	ug/m3
Q1801-01	SS2R	Air	Acetone	6.89		2.28	4.75	ug/m3
Q1801-01	SS2R	Air	Methylene Chloride	3.13	J	2.64	6.95	ug/m3
Q1801-01	SS2R	Air	Toluene	6.78	J	1.66	7.54	ug/m3
Q1801-01	SS2R	Air	Tetrachloroethene	365		0.41	0.81	ug/m3
Q1801-01	SS2R	Air	1,2,4-Trimethylbenzene	2.41	J	1.52	9.83	ug/m3
Q1801-01	SS2R	Air	Hexane	28.9		1.55	7.05	ug/m3
Total Voc :				415				
Total Concentration:				415				
Client ID:	SS3							
Q1801-02	SS3	Air	Acetone	14.0		2.28	4.75	ug/m3
Q1801-02	SS3	Air	2-Butanone	1.53	J	1.42	5.90	ug/m3
Q1801-02	SS3	Air	Chloroform	2.44	J	1.61	9.77	ug/m3
Q1801-02	SS3	Air	Toluene	5.28	J	1.66	7.54	ug/m3
Q1801-02	SS3	Air	Tetrachloroethene	28.5		0.41	0.81	ug/m3
Q1801-02	SS3	Air	m/p-Xylene	4.34	J	3.65	17.4	ug/m3
Q1801-02	SS3	Air	1,2,4-Trimethylbenzene	4.42	J	1.52	9.83	ug/m3
Q1801-02	SS3	Air	Hexane	14.4		1.55	7.05	ug/m3
Total Voc :				75.0				
Total Concentration:				75.0				



QC SUMMARY

Surrogate Summary

SDG No.: Q1801

Client: GFE LLC

Analytical Method: SWTO-15

Lab Sample ID	Client ID	Parameter	Spike	Result	RecoveryQual	Limits	
						Low	High
Q1801-01	SS2R	1-Bromo-4-Fluorobenzene	10	10.5	105	65	135
Q1801-02	SS3	1-Bromo-4-Fluorobenzene	10	10.6	106	65	135
Q1844-02DUP	IA1DUP	1-Bromo-4-Fluorobenzene	10	10.7	107	65	135
VL0421ABL01	VL0421ABL01	1-Bromo-4-Fluorobenzene	10	10.5	105	65	135
VL0421ABS01	VL0421ABS01	1-Bromo-4-Fluorobenzene	10	10.8	108	65	135

Laboratory Control Sample/Laboratory Control Sample Duplicate Summary

SW-846

SDG No.: Q1801

Client: GFE LLC

Analytical Method: SWTO-15

Datafile : VL042397.D

Lab Sample ID	Parameter	Spike	Result	Unit	Rec	RPD	Qual	Low	Limits High	RPD
VL0421ABS01	Dichlorodifluoromethane	10	10.2	ppbv	102			70	130	
	Chloromethane	10	9.60	ppbv	96			70	130	
	Vinyl Chloride	10	8.60	ppbv	86			70	130	
	Bromomethane	10	9.10	ppbv	91			70	130	
	Chloroethane	10	9.50	ppbv	95			70	130	
	Tetrahydrofuran	10	7.80	ppbv	78			70	130	
	Trichlorofluoromethane	10	10.3	ppbv	103			70	130	
	1,1,2-Trichlorotrifluoroethane	10	9.60	ppbv	96			70	130	
	Dichlorotetrafluoroethane	10	10.1	ppbv	101			70	130	
	tert-Butyl Alcohol	10	9.00	ppbv	90			70	130	
	Heptane	10	7.30	ppbv	73			70	130	
	1,1-Dichloroethene	10	9.40	ppbv	94			70	130	
	Acetone	10	8.70	ppbv	87			70	130	
	Carbon disulfide	10	9.40	ppbv	94			70	130	
	Methyl tert-butyl Ether	10	8.50	ppbv	85			70	130	
	Methylene Chloride	10	8.00	ppbv	80			70	130	
	trans-1,2-Dichloroethene	10	9.10	ppbv	91			70	130	
	1,1-Dichloroethane	10	9.60	ppbv	96			70	130	
	Cyclohexane	10	8.50	ppbv	85			70	130	
	2-Butanone	10	8.30	ppbv	83			70	130	
	Carbon Tetrachloride	10	9.50	ppbv	95			70	130	
	cis-1,2-Dichloroethene	10	9.10	ppbv	91			70	130	
	Chloroform	10	9.60	ppbv	96			70	130	
	1,1,1-Trichloroethane	10	9.30	ppbv	93			70	130	
	2,2,4-Trimethylpentane	10	8.10	ppbv	81			70	130	
	Benzene	10	8.60	ppbv	86			70	130	
	1,2-Dichloroethane	10	10.3	ppbv	103			70	130	
	Trichloroethene	10	7.80	ppbv	78			70	130	
	1,2-Dichloropropane	10	8.30	ppbv	83			70	130	
	Bromodichloromethane	10	9.70	ppbv	97			70	130	
	4-Methyl-2-Pentanone	10	7.20	ppbv	72			70	130	
	Toluene	10	8.30	ppbv	83			70	130	
	t-1,3-Dichloropropene	10	8.50	ppbv	85			70	130	
	cis-1,3-Dichloropropene	10	8.30	ppbv	83			70	130	
	1,1,2-Trichloroethane	10	9.00	ppbv	90			70	130	
	Dibromochloromethane	10	9.20	ppbv	92			70	130	
	1,2-Dibromoethane	10	8.90	ppbv	89			70	130	
	Tetrachloroethene	10	7.40	ppbv	74			70	130	
	Chlorobenzene	10	8.90	ppbv	89			70	130	
	Ethyl Benzene	10	9.00	ppbv	90			70	130	
	m/p-Xylene	20	18.2	ppbv	91			70	130	
	o-Xylene	10	9.20	ppbv	92			70	130	
	Styrene	10	8.60	ppbv	86			70	130	
	Bromoform	10	8.80	ppbv	88			70	130	
	1,1,2,2-Tetrachloroethane	10	8.40	ppbv	84			70	130	
	2-Chlorotoluene	10	9.20	ppbv	92			70	130	
	1,3,5-Trimethylbenzene	10	9.00	ppbv	90			70	130	
	1,2,4-Trimethylbenzene	10	8.60	ppbv	86			70	130	

Laboratory Control Sample/Laboratory Control Sample Duplicate Summary

SW-846

SDG No.: Q1801

Client: GFE LLC

Analytical Method: SWTO-15

Datafile : VL042397.D

Lab Sample ID	Parameter	Spike	Result	Unit	Rec	RPD	Qual	Low	Limits High	RPD
VL0421ABS01	1,3-Dichlorobenzene	10	8.70	ppbv	87			70	130	
	1,4-Dichlorobenzene	10	8.60	ppbv	86			70	130	
	1,2-Dichlorobenzene	10	8.70	ppbv	87			70	130	
	1,2,4-Trichlorobenzene	10	8.50	ppbv	85			70	130	
	Hexachloro-1,3-butadiene	10	8.50	ppbv	85			70	130	
	Naphthalene	10	8.20	ppbv	82			70	130	
	1,3-Butadiene	10	9.00	ppbv	90			70	130	
	4-Ethyltoluene	10	9.00	ppbv	90			70	130	
	Hexane	10	7.80	ppbv	78			70	130	
	Allyl Chloride	10	9.10	ppbv	91			70	130	
	1,4-Dioxane	10	7.40	ppbv	74			70	130	
	Methyl methacrylate	10	7.70	ppbv	77			70	130	

Duplicate Sample Summary

Lab Sample Id :	Q1844-02DUP	Q1844-02
Client Id :	IA1DUP	IA1
DF :	1	1
Datafile :	VL042403.D	VL042402.D
Anal Date & Time :	04/21/2025 14:13	04/21/2025 13:38

Parameter	Result	Result	RPD
1,1,1-Trichloroethane	0	0	0
1,1,2,2-Tetrachloroethane	0	0	0
1,1,2-Trichloroethane	0	0	0
1,1,2-Trichlorotrifluoroethane	0	0	0
1,1-Dichloroethane	0	0	0
1,1-Dichloroethene	0	0	0
1,2,4-Trichlorobenzene	0	0	0
1,2,4-Trimethylbenzene	0.23	0.23	0
1,2-Dibromoethane	0	0	0
1,2-Dichlorobenzene	0	0	0
1,2-Dichloroethane	0	0	0
1,2-Dichloropropane	0	0	0
1,3,5-Trimethylbenzene	0	0	0
1,3-Butadiene	0	0	0
1,3-Dichlorobenzene	0	0	0
1,4-Dichlorobenzene	0	0	0
1,4-Dioxane	0	0	0
2,2,4-Trimethylpentane	0.35	0.32	9
2-Butanone	0.66	0.68	3
2-Chlorotoluene	0	0	0
4-Ethyltoluene	0	0	0
4-Methyl-2-Pentanone	0	0	0
Acetone	22.8	22.7	0.44
Allyl Chloride	0	0	0
Benzene	12.8	12.5	2.4
Bromodichloromethane	0	0	0
Bromoform	0	0	0
Bromomethane	0	0	0
Carbon Disulfide	0	0	0
Carbon Tetrachloride	0.07	0.07	0

Duplicate Sample Summary

Lab Sample Id :	Q1844-02DUP	Q1844-02
Client Id :	IA1DUP	IA1
DF :	1	1
Datafile :	VL042403.D	VL042402.D
Anal Date & Time :	04/21/2025 14:13	04/21/2025 13:38

Parameter	Result	Result	RPD
Chlorobenzene	0	0	0
Chloroethane	0	0	0
Chloroform	0	0	0
Chloromethane	0.51	0.5	2
cis-1,2-Dichloroethene	0	0	0
cis-1,3-Dichloropropene	0	0	0
Cyclohexane	0.22	0	200 *
Dibromochloromethane	0	0	0
Dichlorodifluoromethane	0.39	0.41	5
Dichlorotetrafluoroethane	0	0	0
Ethyl Benzene	0.29	0.28	3.5
Heptane	0.37	0.35	5.6
Hexachloro-1,3-Butadiene	0	0	0
Hexane	0.72	0.67	7.2
m/p-Xylene	0.82	0.79	3.7
Methyl Methacrylate	0	0	0
Methyl tert-Butyl Ether	0	0	0
Methylene Chloride	0.62	0.6	3.3
Naphthalene	0.51	0.52	1.9
o-Xylene	0.32	0.32	0
Styrene	0.39	0.36	8
t-1,3-Dichloropropene	0	0	0
tert-Butyl alcohol	5.1	5.1	0
Tetrachloroethene	19.5	19.3	1
Tetrahydrofuran	0	0	0
Toluene	2	2	0
trans-1,2-Dichloroethene	0	0	0
Trichloroethene	0	0	0
Trichlorofluoromethane	0.24	0.23	4.3
Vinyl Chloride	0	0	0

VOLATILE METHOD BLANK SUMMARY

EPA SAMPLE NO.

VL0421ABL01

Lab Name: CHEMTECH

Contract: GFEL01

Lab Code: CHEM Case No.: Q1801

SAS No.: Q1801 SDG NO.: Q1801

Lab File ID: VL042396.D

Lab Sample ID: VL0421ABL01

Date Analyzed: 04/21/2025

Time Analyzed: 09:28

GC Column: RTX-1 ID: 0.32 (mm)

Heated Purge: (Y/N) N

Instrument ID: MSVOA_L

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES, MS AND MSD:

EPA SAMPLE NO.	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED
VL0421ABS01	VL0421ABS01	VL042397.D	04/21/2025
SS2R	Q1801-01	VL042400.D	04/21/2025
SS3	Q1801-02	VL042401.D	04/21/2025
IA1DUP	Q1844-02DUP	VL042403.D	04/21/2025

COMMENTS: _____

VOLATILE ORGANIC INSTRUMENT PERFORMANCE CHECK
BROMOFLUOROBENZENE (BFB)

Lab Name: CHEMTECH Contract: GFEL01
Lab Code: CHEM Case No.: Q1801 SAS No.: Q1801 SDG NO.: Q1801
Lab File ID: VL042348.D BFB Injection Date: 04/17/2025
Instrument ID: MSVOA_L BFB Injection Time: 07:50
GC Column: RTX-1 ID: 0.32 (mm) Heated Purge: Y/N N

m/e	ION ABUNDANCE CRITERIA	% RELATIVE ABUNDANCE
50	8.0 - 40.0% of mass 95	23.1
75	30.0 - 66.0% of mass 95	57
95	Base Peak, 100% relative abundance	100
96	5.0 - 9.0% of mass 95	6.8
173	Less than 2.0% of mass 174	0.0 (0.0) 1
174	50.0 - 120.0% of mass 95	59.5
175	4.0 - 9.0% of mass 174	4.5 (7.6) 1
176	93.0 - 101.0% of mass 174	57.6 (96.9) 1
177	5.0 - 9.0% of mass 176	3.6 (6.3) 2

1-Value is % mass 174

2-Value is % mass 176

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS, AND STANDARDS:

EPA SAMPLE NO.	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED	TIME ANALYZED
VSTDICCC010	VSTDICCC010	VL042349.D	04/17/2025	09:44
VSTDICCC002	VSTDICCC002	VL042350.D	04/17/2025	10:17
VSTDICCC001	VSTDICCC001	VL042351.D	04/17/2025	10:48
VSTDICCC0.5	VSTDICCC0.5	VL042352.D	04/17/2025	11:19
VSTDICCC0.1	VSTDICCC0.1	VL042353.D	04/17/2025	11:50
VSTDICCC0.03	VSTDICCC0.03	VL042354.D	04/17/2025	12:22
VSTDICCC015	VSTDICCC015	VL042355.D	04/17/2025	12:54

VOLATILE ORGANIC INSTRUMENT PERFORMANCE CHECK
BROMOFLUOROBENZENE (BFB)

Lab Name: CHEMTECH Contract: GFEL01
Lab Code: CHEM Case No.: Q1801 SAS No.: Q1801 SDG NO.: Q1801
Lab File ID: VL042394.D BFB Injection Date: 04/21/2025
Instrument ID: MSVOA_L BFB Injection Time: 07:58
GC Column: RTX-1 ID: 0.32 (mm) Heated Purge: Y/N N

m/e	ION ABUNDANCE CRITERIA	% RELATIVE ABUNDANCE
50	8.0 - 40.0% of mass 95	26.7
75	30.0 - 66.0% of mass 95	63.9
95	Base Peak, 100% relative abundance	100
96	5.0 - 9.0% of mass 95	7
173	Less than 2.0% of mass 174	0.0 (0.0) 1
174	50.0 - 120.0% of mass 95	54.7
175	4.0 - 9.0% of mass 174	4.2 (7.7) 1
176	93.0 - 101.0% of mass 174	52.9 (96.6) 1
177	5.0 - 9.0% of mass 176	3.4 (6.4) 2

1-Value is % mass 174

2-Value is % mass 176

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS, AND STANDARDS:

EPA SAMPLE NO.	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED	TIME ANALYZED
VSTDCCC010	VSTDCCC010	VL042395.D	04/21/2025	08:41
VL0421ABL01	VL0421ABL01	VL042396.D	04/21/2025	09:28
VL0421ABS01	VL0421ABS01	VL042397.D	04/21/2025	10:14
SS2R	Q1801-01	VL042400.D	04/21/2025	12:07
SS3	Q1801-02	VL042401.D	04/21/2025	12:51
IA1DUP	Q1844-02DUP	VL042403.D	04/21/2025	14:13

VOLATILE INTERNAL STANDARD AREA AND RT SUMMARY

Lab Name: CHEMTECH Contract: GFEL01
 Lab Code: CHEM Case No.: Q1801 SAS No.: Q1801 SDG NO.: Q1801
 Lab File ID: VL042395.D Date Analyzed: 04/21/2025
 Instrument ID: MSVOA_L Time Analyzed: 08:41
 GC Column: RTX-1 ID: 0.32 (mm) Heated Purge: (Y/N) N

	IS1 AREA #	RT #	IS2 AREA #	RT #	IS3 AREA #	RT #
12 HOUR STD	118349	2.78	334726	3.95	297161	8.88
UPPER LIMIT	165689	3.11	468616	4.28	416025	9.21
LOWER LIMIT	71009.4	2.45	200836	3.62	178297	8.55
EPA SAMPLE NO.						
SS2R	127438	2.78	352813	3.95	303522	8.88
SS2R	127438	2.78	352813	3.95	303522	8.88
SS3	122041	2.79	337832	3.96	286610	8.88
SS3	122041	2.79	337832	3.96	286610	8.88
IA1DUP	132791	2.78	375368	3.95	317335	8.88
IA1DUP	132791	2.78	375368	3.95	317335	8.88
VL0421ABL01	120411	2.78	341628	3.95	293743	8.88
VL0421ABL01	120411	2.78	341628	3.95	293743	8.88
VL0421ABS01	118401	2.78	337921	3.95	293537	8.88
VL0421ABS01	118401	2.78	337921	3.95	293537	8.88

IS1 = Bromochloromethane
 IS2 = 1,4-Difluorobenzene
 IS3 = Chlorobenzene-d5

AREA UPPER LIMIT = +40% of internal standard area
 AREA LOWER LIMIT = -40% of internal standard area
 RT UPPER LIMIT = +0.33 minutes of internal standard RT
 RT LOWER LIMIT = -0.33 minutes of internal standard RT

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



SAMPLE DATA

Report of Analysis

Client:	GFE LLC	Date Collected:	04/12/25
Project:	416 Clinton St Hempstead	Date Received:	04/14/25
Client Sample ID:	SS2R	SDG No.:	Q1801
Lab Sample ID:	Q1801-01	Matrix:	Air
Analytical Method:	TO-15	Test:	TO-15
Sample Wt/Vol:	400	Units:	mL

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
VL042400.D	4		04/21/25 12:07	VL042125

CAS Number	Parameter	Conc. ppbv	Conc. ug/M3	Qualifier	MDL	LOQ / CRQL	Units
TARGETS							
75-71-8	Dichlorodifluoromethane	0.84	4.15	U	4.15	9.89	ug/m3
74-87-3	Chloromethane	0.39	0.81	U	0.81	4.13	ug/m3
75-01-4	Vinyl Chloride	0.060	0.15	U	0.15	0.31	ug/m3
74-83-9	Bromomethane	0.56	2.17	U	2.17	7.77	ug/m3
75-00-3	Chloroethane	0.68	1.79	U	1.79	5.28	ug/m3
109-99-9	Tetrahydrofuran	0.52	1.53	U	1.53	5.90	ug/m3
75-69-4	Trichlorofluoromethane	0.64	3.60	U	3.60	11.2	ug/m3
76-13-1	1,1,2-Trichlorotrifluoroethane	0.64	4.91	U	4.91	15.3	ug/m3
76-14-2	Dichlorotetrafluoroethane	0.35	2.45	U	2.45	14.0	ug/m3
75-65-0	tert-Butyl alcohol	0.76	2.30	U	2.30	6.06	ug/m3
142-82-5	Heptane	0.46	1.89	J	1.80	8.20	ug/m3
75-35-4	1,1-Dichloroethene	0.56	2.22	U	2.22	7.93	ug/m3
67-64-1	Acetone	2.90	6.89		2.28	4.75	ug/m3
75-15-0	Carbon Disulfide	0.64	1.99	U	1.99	6.23	ug/m3
1634-04-4	Methyl tert-Butyl Ether	0.48	1.73	U	1.73	7.21	ug/m3
75-09-2	Methylene Chloride	0.90	3.13	J	2.64	6.95	ug/m3
156-60-5	trans-1,2-Dichloroethene	0.76	3.01	U	3.01	7.93	ug/m3
75-34-3	1,1-Dichloroethane	0.52	2.10	U	2.10	8.09	ug/m3
110-82-7	Cyclohexane	0.88	3.03	U	3.03	6.88	ug/m3
78-93-3	2-Butanone	0.48	1.42	U	1.42	5.90	ug/m3
56-23-5	Carbon Tetrachloride	0.040	0.25	U	0.25	0.75	ug/m3
156-59-2	cis-1,2-Dichloroethene	0.36	1.43	U	1.43	7.93	ug/m3
67-66-3	Chloroform	0.33	1.61	U	1.61	9.77	ug/m3
71-55-6	1,1,1-Trichloroethane	0.040	0.22	U	0.22	0.65	ug/m3
540-84-1	2,2,4-Trimethylpentane	0.40	1.87	U	1.87	9.34	ug/m3
71-43-2	Benzene	0.35	1.12	U	1.12	6.39	ug/m3
107-06-2	1,2-Dichloroethane	0.36	1.46	U	1.46	8.09	ug/m3
79-01-6	Trichloroethene	0.070	0.38	U	0.38	0.64	ug/m3
78-87-5	1,2-Dichloropropane	0.44	2.03	U	2.03	9.24	ug/m3
75-27-4	Bromodichloromethane	0.31	2.08	U	2.08	13.4	ug/m3
108-10-1	4-Methyl-2-Pentanone	0.37	1.52	U	1.52	8.20	ug/m3
108-88-3	Toluene	1.80	6.78	J	1.66	7.54	ug/m3
10061-02-6	t-1,3-Dichloropropene	0.24	1.09	U	1.09	9.08	ug/m3
10061-01-5	cis-1,3-Dichloropropene	0.26	1.18	U	1.18	9.08	ug/m3
79-00-5	1,1,2-Trichloroethane	0.28	1.53	U	1.53	10.9	ug/m3

Report of Analysis

Client:	GFE LLC	Date Collected:	04/12/25
Project:	416 Clinton St Hempstead	Date Received:	04/14/25
Client Sample ID:	SS2R	SDG No.:	Q1801
Lab Sample ID:	Q1801-01	Matrix:	Air
Analytical Method:	TO-15	Test:	TO-15
Sample Wt/Vol:	400	Units:	mL

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
VL042400.D	4		04/21/25 12:07	VL042125

CAS Number	Parameter	Conc. ppbv	Conc. ug/M3	Qualifier	MDL	LOQ / CRQL	Units
124-48-1	Dibromochloromethane	0.26	2.22	U	2.22	17.0	ug/m3
106-93-4	1,2-Dibromoethane	0.29	2.23	U	2.23	3.07	ug/m3
127-18-4	Tetrachloroethene	53.8	365		0.41	0.81	ug/m3
108-90-7	Chlorobenzene	0.30	1.38	U	1.38	9.21	ug/m3
100-41-4	Ethyl Benzene	0.48	2.08	U	2.08	8.69	ug/m3
179601-23-1	m/p-Xylene	0.84	3.65	U	3.65	17.4	ug/m3
95-47-6	o-Xylene	0.48	2.08	U	2.08	8.69	ug/m3
100-42-5	Styrene	0.44	1.87	U	1.87	8.52	ug/m3
75-25-2	Bromoform	0.23	2.38	U	2.38	20.7	ug/m3
79-34-5	1,1,2,2-Tetrachloroethane	0.040	0.27	U	0.27	0.82	ug/m3
95-49-8	2-Chlorotoluene	0.44	2.28	U	2.28	10.4	ug/m3
108-67-8	1,3,5-Trimethylbenzene	0.44	2.16	U	2.16	9.83	ug/m3
95-63-6	1,2,4-Trimethylbenzene	0.49	2.41	J	1.52	9.83	ug/m3
541-73-1	1,3-Dichlorobenzene	0.32	1.92	U	1.92	12.0	ug/m3
106-46-7	1,4-Dichlorobenzene	0.21	1.26	U	1.26	12.0	ug/m3
95-50-1	1,2-Dichlorobenzene	0.30	1.80	U	1.80	12.0	ug/m3
120-82-1	1,2,4-Trichlorobenzene	0.34	2.52	U	2.52	14.8	ug/m3
87-68-3	Hexachloro-1,3-Butadiene	0.34	3.63	U	3.63	21.3	ug/m3
106-99-0	1,3-Butadiene	0.52	1.15	U	1.15	4.42	ug/m3
91-20-3	Naphthalene	0.30	1.57	U	1.57	2.10	ug/m3
622-96-8	4-Ethyltoluene	0.48	2.36	U	2.36	9.83	ug/m3
110-54-3	Hexane	8.20	28.9		1.55	7.05	ug/m3
107-05-1	Allyl Chloride	0.60	1.88	U	1.88	6.26	ug/m3
123-91-1	1,4-Dioxane	0.84	3.03	U	3.03	7.21	ug/m3
80-62-6	Methyl Methacrylate	0.34	1.39	U	1.39	8.19	ug/m3
SURROGATES							
460-00-4	1-Bromo-4-Fluorobenzene	10.5			65 - 135	105%	SPK: 10
INTERNAL STANDARDS							
74-97-5	Bromochloromethane	127000		2.784			
540-36-3	1,4-Difluorobenzene	353000		3.952			
3114-55-4	Chlorobenzene-d5	304000		8.875			

Report of Analysis

Client:	GFE LLC	Date Collected:	04/12/25
Project:	416 Clinton St Hempstead	Date Received:	04/14/25
Client Sample ID:	SS2R	SDG No.:	Q1801
Lab Sample ID:	Q1801-01	Matrix:	Air
Analytical Method:	TO-15	Test:	TO-15
Sample Wt/Vol:	400	Units:	mL

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
VL042400.D	4		04/21/25 12:07	VL042125

CAS Number	Parameter	Conc. ppbv	Conc. ug/M3	Qualifier	MDL	LOQ / CRQL	Units
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U = Not Detected
 RL = Reporting Limit
 MDL = Method Detection Limit
 E = Value Exceeds Calibration Range
 D = Dilution

J = Estimated Value
 B = Analyte Found in Associated Method Blank
 N = Presumptive Evidence of a Compound
 * = Values outside of QC limits
 Q = indicates LCS control criteria did not meet requirements

Data Path : Z:\voasrv\HPCHEM1\MSVOA_L\Data\VL042125\
 Data File : VL042400.D
 Acq On : 21 Apr 2025 12:07
 Operator : SY/MD
 Sample : Q1801-01 4X
 Misc : 400mL/MSVOA_L
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_L
 ClientSampleId :
 SS2R

Manual Integrations
 APPROVED

Reviewed By :Semsettin Yesilyurt 04/22/2025
 Supervised By :Mahesh Dadoda 04/22/2025

Quant Time: Apr 22 02:52:43 2025

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_L\methods\VL041725AIR.M

Quant Title : AIR ANALYSIS BY METHOD TO-15 Instrument: MSVOA_L Fri Aug 2

QLast Update : Fri Apr 18 02:14:00 2025

Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) Bromochloromethane	2.784	49	127438	10.000	ppbv	0.00
33) 1,4-Difluorobenzene	3.952	114	352813	10.000	ppbv	0.00
55) Chlorobenzene-d5	8.875	117	303522	10.000	ppbv	0.00
System Monitoring Compounds						
68) 1-Bromo-4-Fluorobenzene	10.370	95	261136	10.470	ppbv	0.00
Spiked Amount	10.000	Range	65 - 135	Recovery	=	104.700%
Target Compounds						
					Qvalue	
2) Dichlorodifluoromethane	1.496	85	2752	0.122	ppbv	90
10) Heptane	4.965	43	3144m	0.115	ppbv	
13) Ethanol	1.719	45	2925m	1.885	ppbv	
15) Acetone	1.835	43	15004	0.719	ppbv	93
20) Methylene Chloride	2.062	84	1839	0.225	ppbv #	87
23) Vinyl Acetate	2.460	43	53575	6.344	ppbv #	96
26) Hexane	2.823	57	42883	2.060	ppbv #	26
34) 2-Butanone	2.560	43	2937	0.105	ppbv	98
52) Tetrachloroethene	8.218	164	197775	13.446	ppbv #	86
53) Toluene	6.923	91	20734	0.454	ppbv	96
59) m/p-Xylene	9.526	91	5667m	0.117	ppbv	
74) 1,2,4-Trimethylbenzene	11.529	105	7515	0.122	ppbv #	61

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

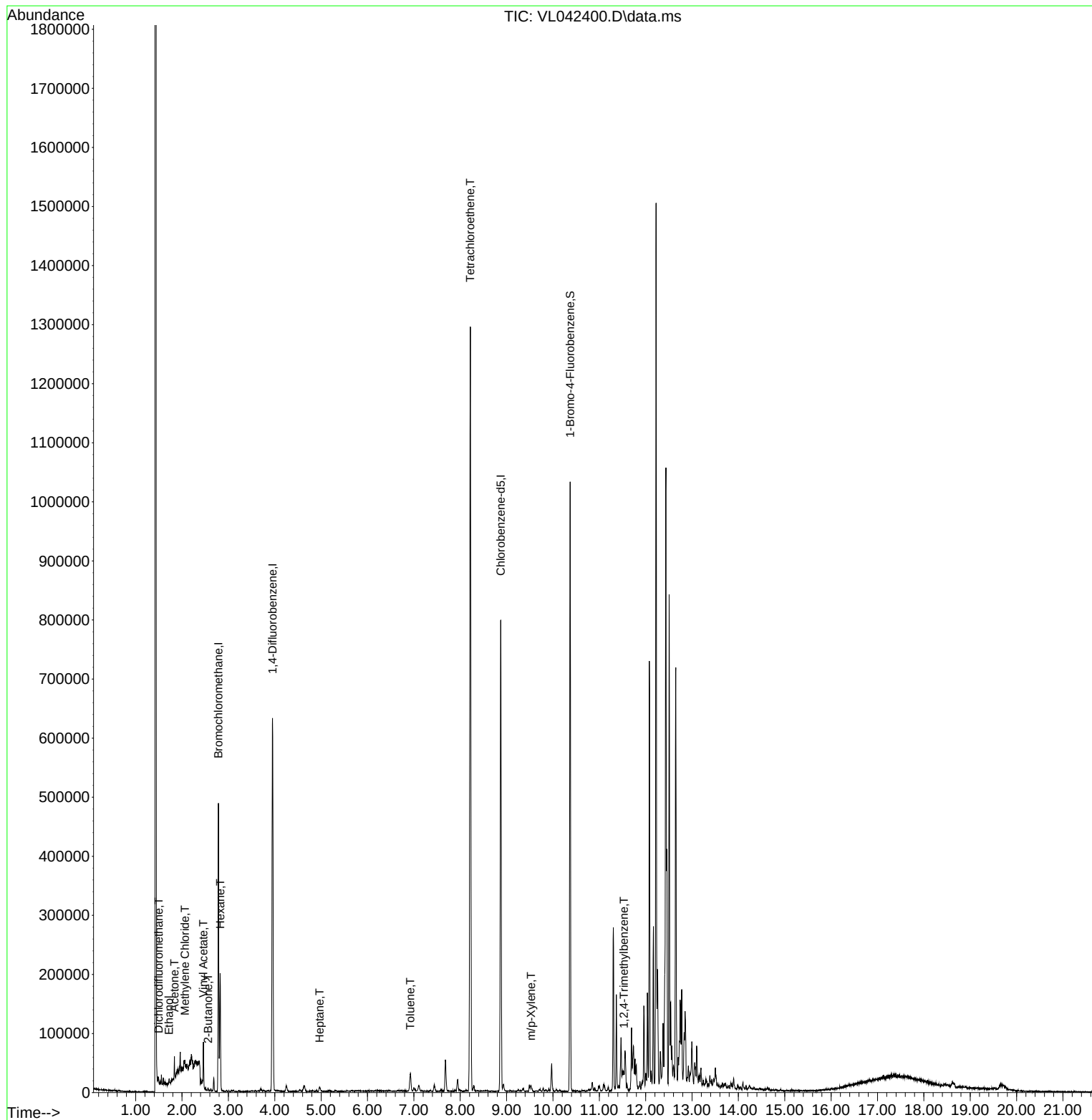
Data Path : Z:\voasrv\HPCHEM1\MSVOA_L\Data\VL042125\
Data File : VL042400.D
Acq On : 21 Apr 2025 12:07
Operator : SY/MD
Sample : Q1801-01 4X
Misc : 400mL/MSVOA_L
ALS Vial : 1 Sample Multiplier: 1

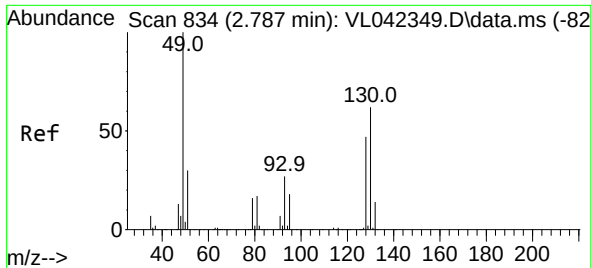
Instrument :
MSVOA_L
ClientSampleId :
SS2R

Quant Time: Apr 22 02:52:43 2025
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_L\methods\VL041725AIR.M
Quant Title : AIR ANALYSIS BY METHOD TO-15 Instrument: MSVOA_L Fri Aug
QLast Update : Fri Apr 18 02:14:00 2025
Response via : Initial Calibration

Manual Integrations
APPROVED

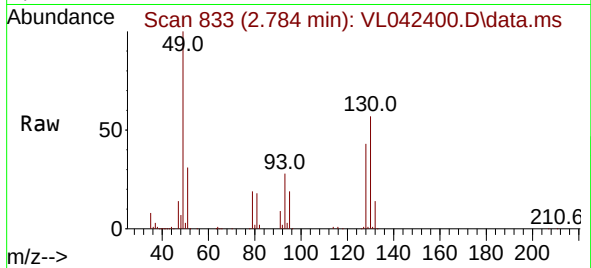
Reviewed By :Semsettin Yesilyurt 04/22/2025
Supervised By :Mahesh Dadoda 04/22/2025





#1
Bromochloromethane
Concen: 10.000 ppbv
RT: 2.784 min Scan# 81
Delta R.T. -0.003 min
Lab File: VL042400.D
Acq: 21 Apr 2025 12:07

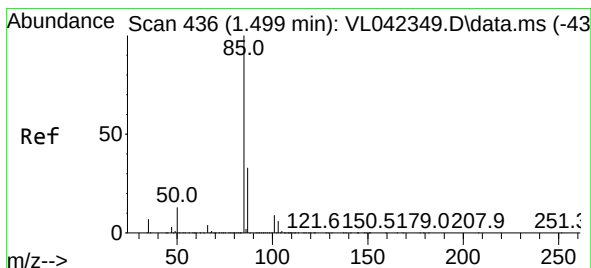
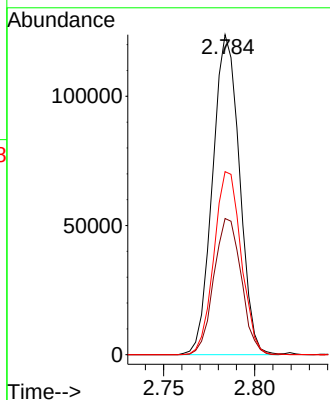
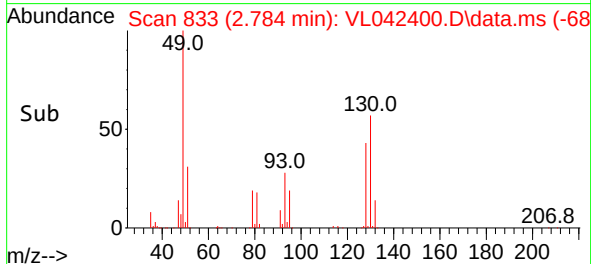
Instrument :
MSVOA_L
ClientSampleId :
SS2R



Tgt Ion: 49 Resp: 127438
Ion Ratio Lower Upper
49 100
128 43.3 22.8 68.4
130 57.8 29.3 88.0

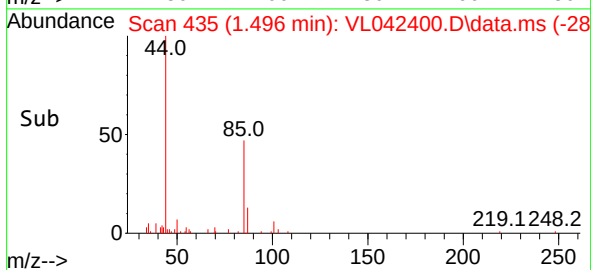
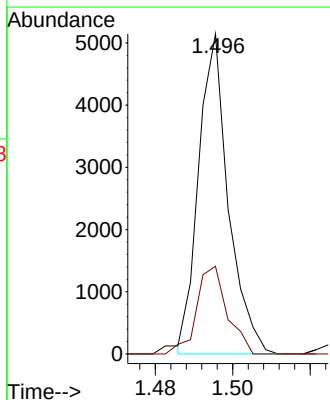
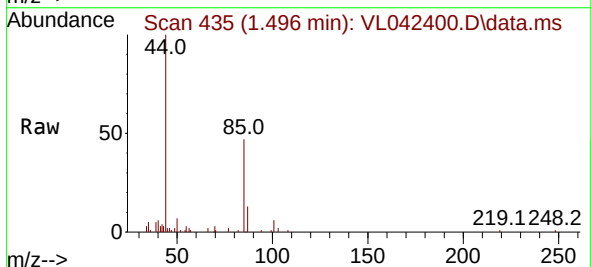
Manual Integrations
APPROVED

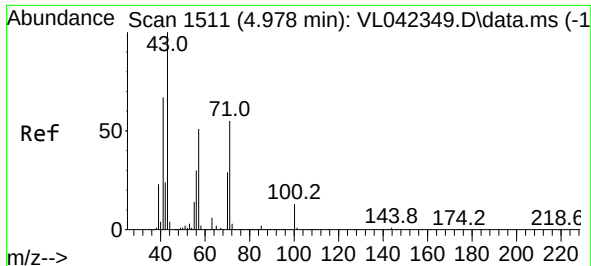
Reviewed By :Semsettin Yesilyurt 04/22/2025
Supervised By :Mahesh Dadoda 04/22/2025



#2
Dichlorodifluoromethane
Concen: 0.122 ppbv
RT: 1.496 min Scan# 435
Delta R.T. -0.003 min
Lab File: VL042400.D
Acq: 21 Apr 2025 12:07

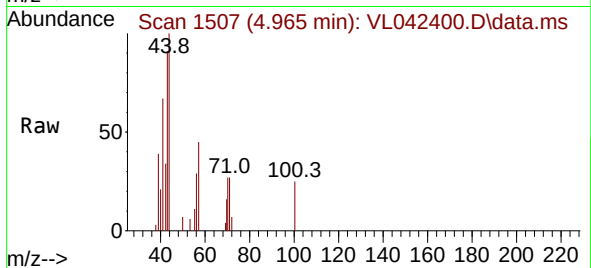
Tgt Ion: 85 Resp: 2752
Ion Ratio Lower Upper
85 100
87 27.4 26.6 39.8





#10
Heptane
Concen: 0.115 ppbv m
RT: 4.965 min Scan# 11
Delta R.T. -0.013 min
Lab File: VL042400.D
Acq: 21 Apr 2025 12:07

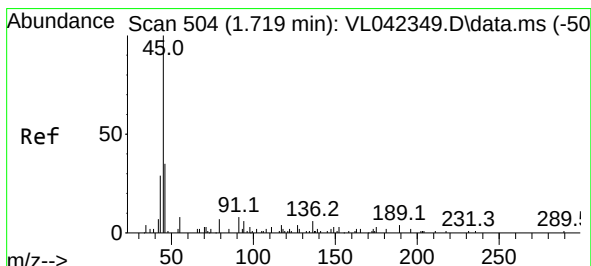
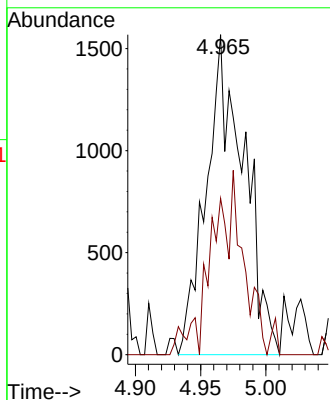
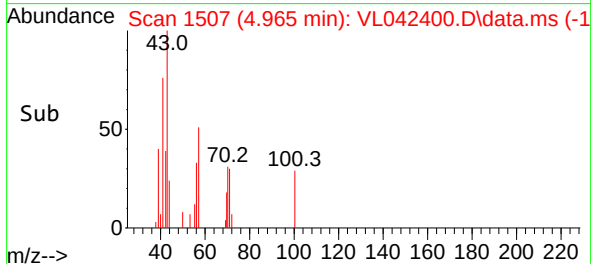
Instrument :
MSVOA_L
ClientSampleId :
SS2R



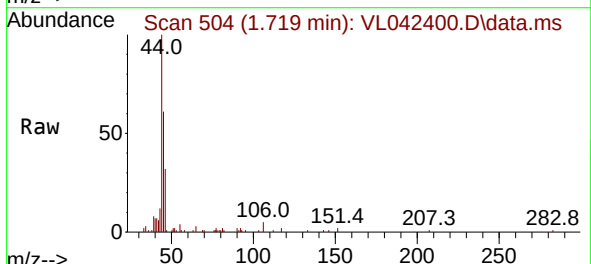
Tgt Ion: 43 Resp: 314
Ion Ratio Lower Upper
43 100
57 50.2 39.9 59.9

Manual Integrations
APPROVED

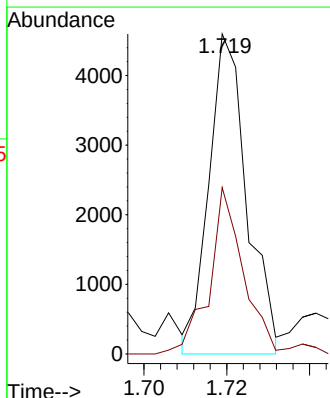
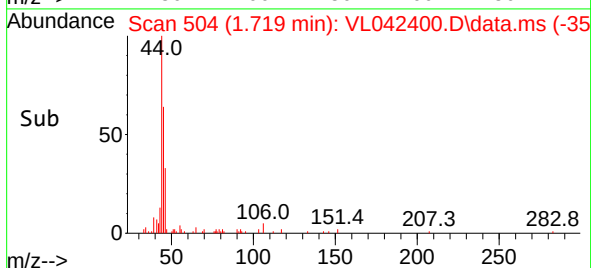
Reviewed By :Semsettin Yesilyurt 04/22/2025
Supervised By :Mahesh Dadoda 04/22/2025

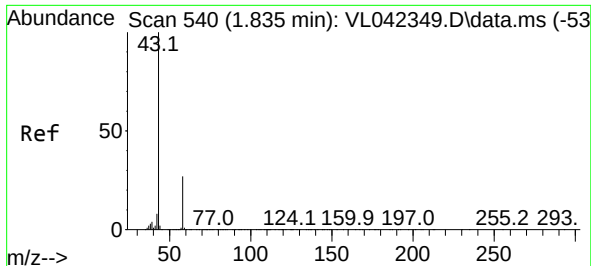


#13
Ethanol
Concen: 1.885 ppbv m
RT: 1.719 min Scan# 504
Delta R.T. 0.000 min
Lab File: VL042400.D
Acq: 21 Apr 2025 12:07



Tgt Ion: 45 Resp: 2925
Ion Ratio Lower Upper
45 100
46 8881.7 21.9 87.7#





#15

Acetone

Concen: 0.719 ppbv

RT: 1.835 min Scan# 540

Delta R.T. 0.000 min

Lab File: VL042400.D

Acq: 21 Apr 2025 12:07

Instrument :

MSVOA_L

ClientSampleId :

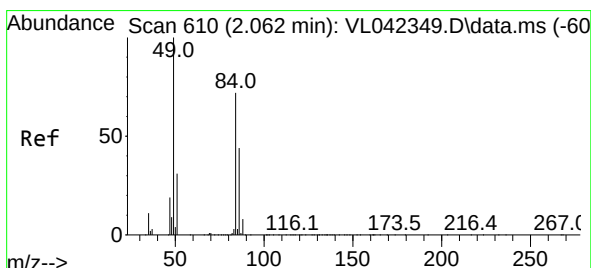
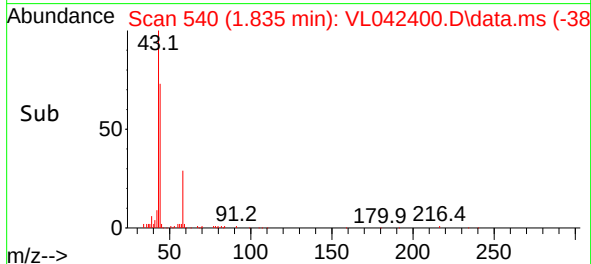
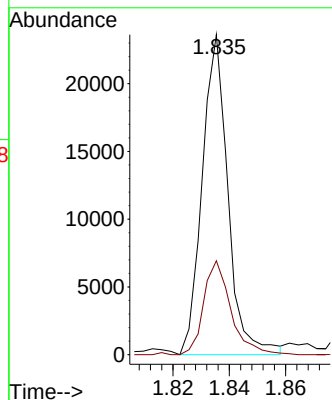
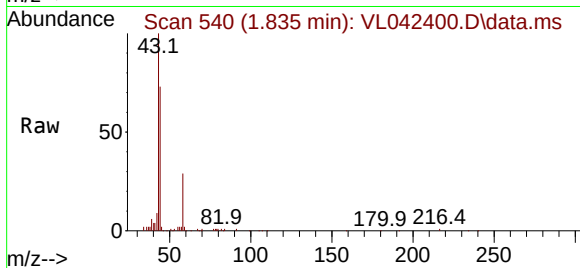
SS2R

Tgt Ion: 43 Resp: 15004
 Ion Ratio Lower Upper
 43 100
 58 31.5 22.4 33.6

Manual Integrations

APPROVED

Reviewed By :Semsettin Yesilyurt 04/22/2025
 Supervised By :Mahesh Dadoda 04/22/2025



#20

Methylene Chloride

Concen: 0.225 ppbv

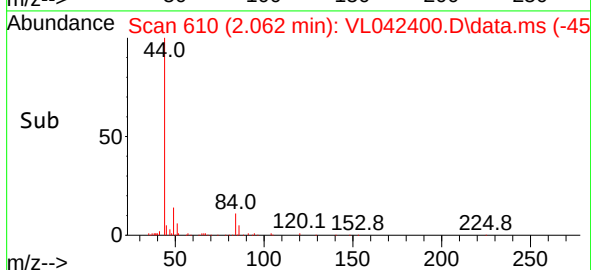
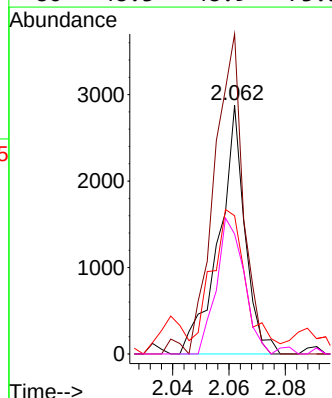
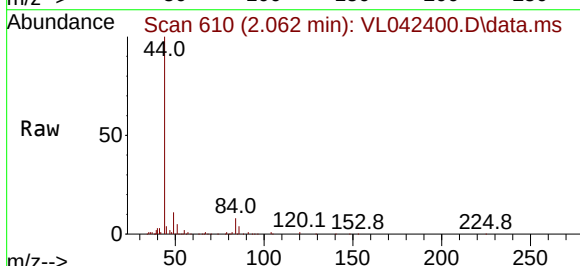
RT: 2.062 min Scan# 610

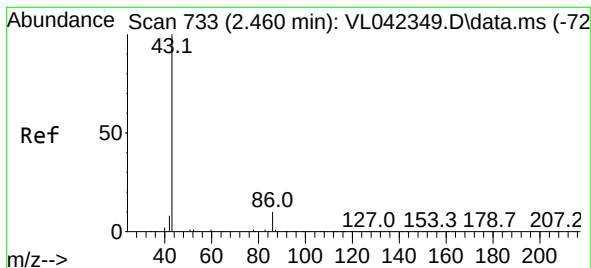
Delta R.T. 0.000 min

Lab File: VL042400.D

Acq: 21 Apr 2025 12:07

Tgt Ion: 84 Resp: 1839
 Ion Ratio Lower Upper
 84 100
 49 128.6 111.6 167.4
 51 56.5 35.1 52.7#
 86 48.3 48.9 73.3#





#23

Vinyl Acetate

Concen: 6.344 ppbv

RT: 2.460 min Scan# 71

Delta R.T. 0.000 min

Lab File: VL042400.D

Acq: 21 Apr 2025 12:07

Instrument :

MSVOA_L

ClientSampleId :

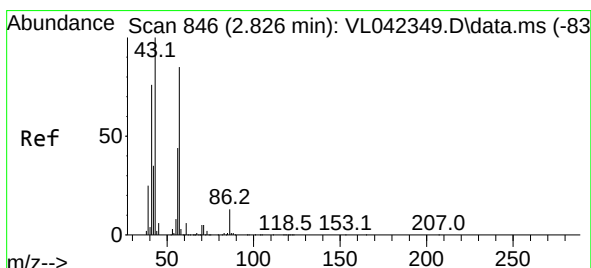
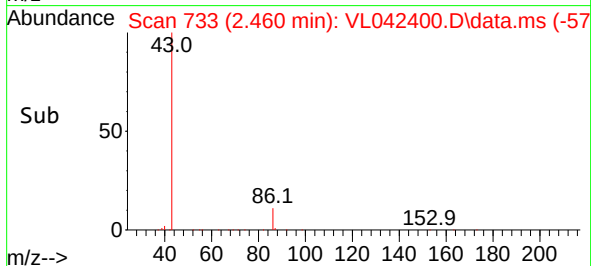
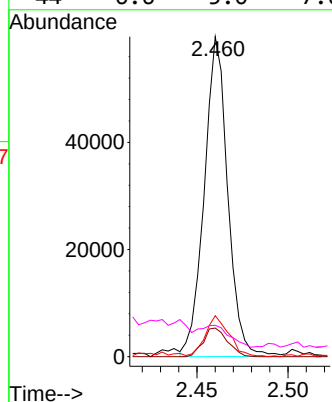
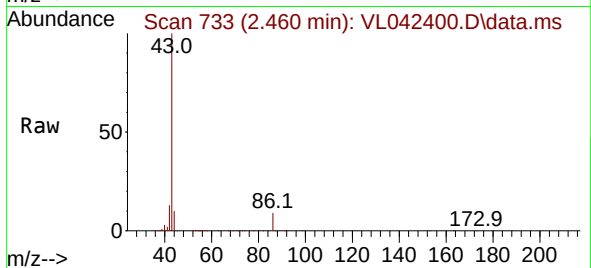
SS2R

Tgt Ion: 43	Resp: 5357
Ion Ratio	Lower Upper
43 100	
86 9.0	6.5 9.7
42 12.9	7.8 11.8
44 6.6	5.0 7.6

Manual Integrations

APPROVED

Reviewed By :Semsettin Yesilyurt 04/22/2025
Supervised By :Mahesh Dadoda 04/22/2025



#26

Hexane

Concen: 2.060 ppbv

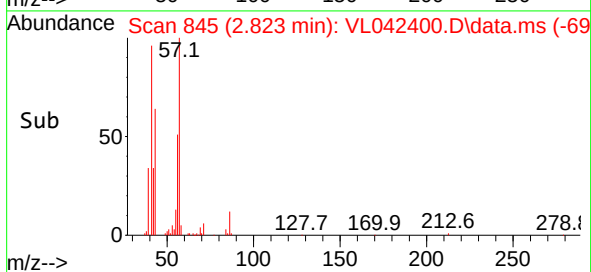
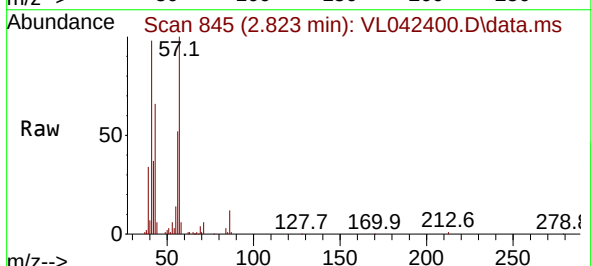
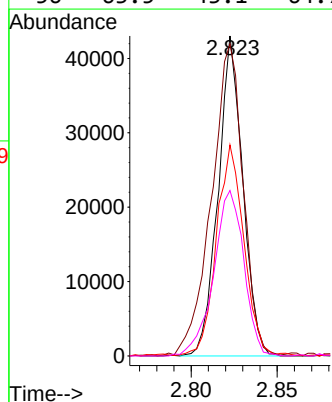
RT: 2.823 min Scan# 845

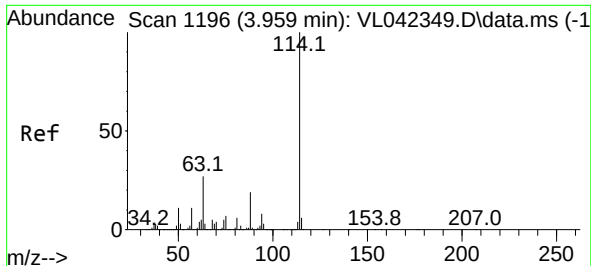
Delta R.T. -0.003 min

Lab File: VL042400.D

Acq: 21 Apr 2025 12:07

Tgt	Ion: 57	Resp:	42883
Ion	Ratio	Lower	Upper
57	100		
41	126.1	77.9	116.9#
43	74.8	208.4	312.6#
56	63.5	43.1	64.7





#33

1,4-Difluorobenzene

Concen: 10.000 ppbv

RT: 3.952 min Scan# 1194

Delta R.T. -0.006 min

Lab File: VL042400.D

Acq: 21 Apr 2025 12:07

Instrument :

MSVOA_L

ClientSampleId :

SS2R

Tgt Ion:114 Resp: 35281

Ion Ratio Lower Upper

114 100

63 28.1 21.8 32.6

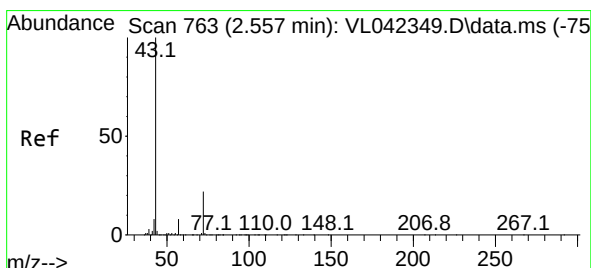
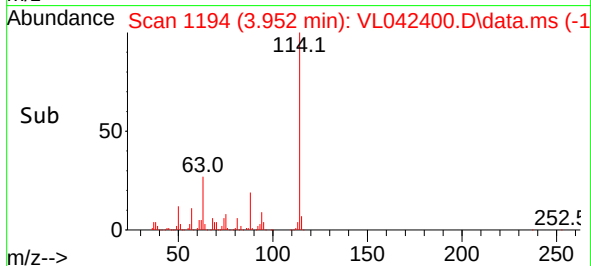
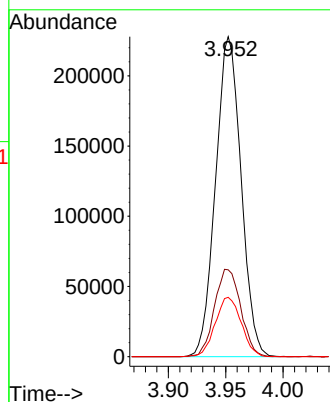
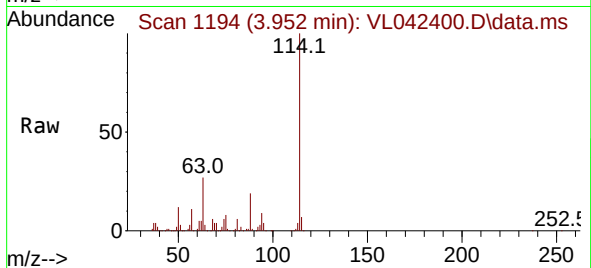
88 18.9 15.5 23.3

Manual Integrations

APPROVED

Reviewed By :Semsettin Yesilyurt 04/22/2025

Supervised By :Mahesh Dadoda 04/22/2025



#34

2-Butanone

Concen: 0.105 ppbv

RT: 2.560 min Scan# 764

Delta R.T. 0.003 min

Lab File: VL042400.D

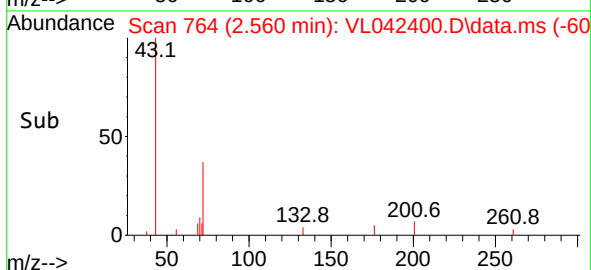
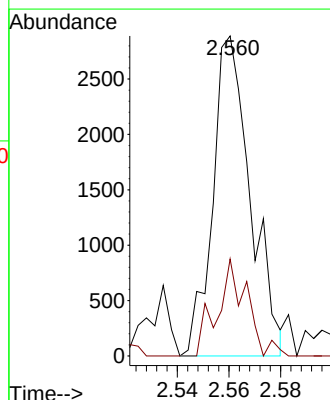
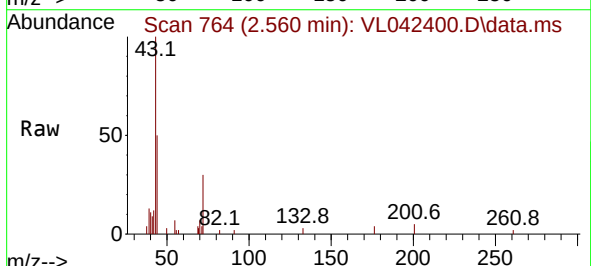
Acq: 21 Apr 2025 12:07

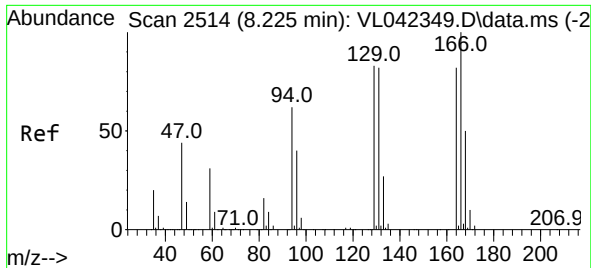
Tgt Ion: 43 Resp: 2937

Ion Ratio Lower Upper

43 100

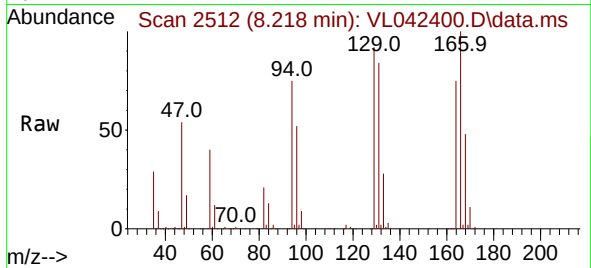
72 23.9 18.2 27.2





#52
Tetrachloroethene
Concen: 13.446 ppbv
RT: 8.218 min Scan# 2112
Delta R.T. -0.006 min
Lab File: VL042400.D
Acq: 21 Apr 2025 12:07

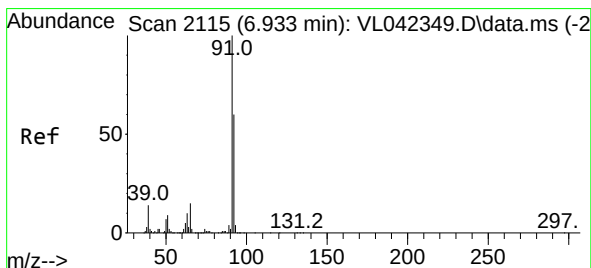
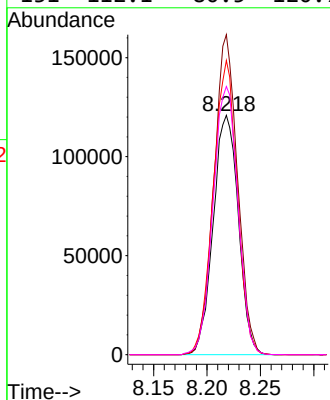
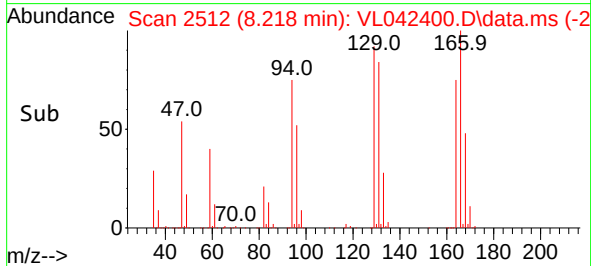
Instrument :
MSVOA_L
ClientSampleId :
SS2R



Tgt Ion:164 Resp: 197775
Ion Ratio Lower Upper
164 100
166 133.7 98.2 147.2
129 123.0 81.8 122.8
131 112.1 80.5 120.7

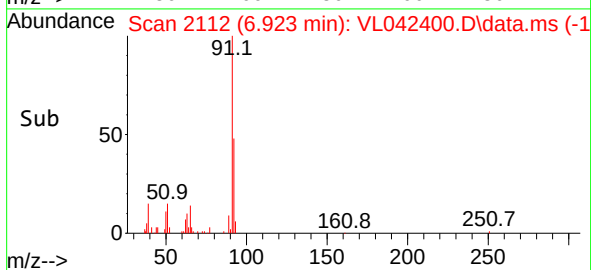
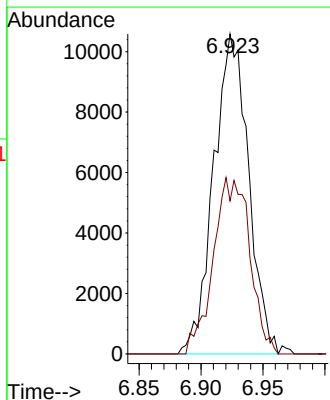
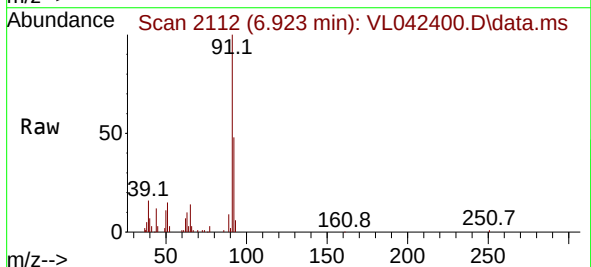
Manual Integrations
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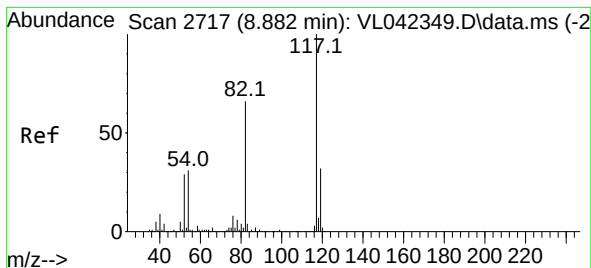
Reviewed By :Semsettin Yesilyurt 04/22/2025
Supervised By :Mahesh Dadoda 04/22/2025



#53
Toluene
Concen: 0.454 ppbv
RT: 6.923 min Scan# 2112
Delta R.T. -0.010 min
Lab File: VL042400.D
Acq: 21 Apr 2025 12:07

Tgt Ion: 91 Resp: 20734
Ion Ratio Lower Upper
91 100
92 57.6 48.2 72.4





#55

Chlorobenzene-d5

Concen: 10.000 ppbv

RT: 8.875 min Scan# 21

Delta R.T. -0.006 min

Lab File: VL042400.D

Acq: 21 Apr 2025 12:07

Instrument :

MSVOA_L

ClientSampleId :

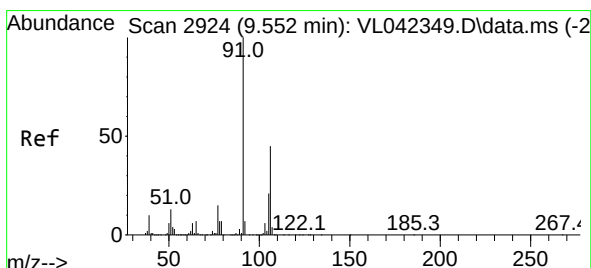
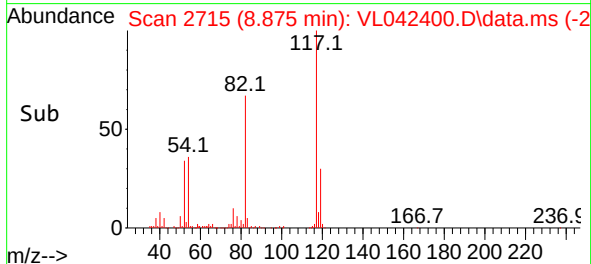
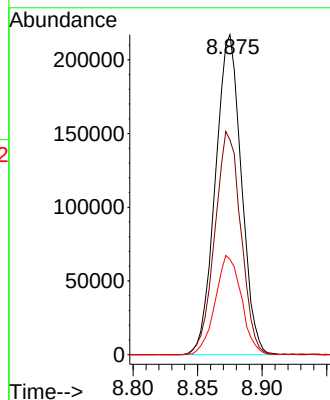
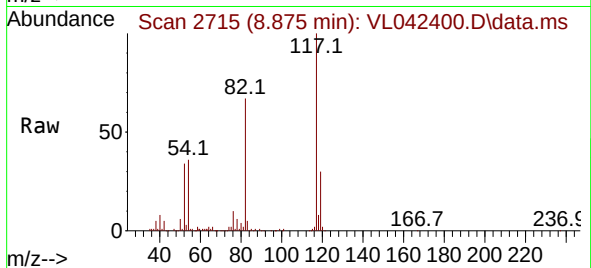
SS2R

Tgt Ion	Ratio	Lower	Upper
117	100		
82	70.2	57.8	86.6
119	31.3	25.2	37.8

Manual Integrations

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Reviewed By :Semsettin Yesilyurt 04/22/2025
Supervised By :Mahesh Dadoda 04/22/2025



#59

m/p-Xylene

Concen: 0.117 ppbv m

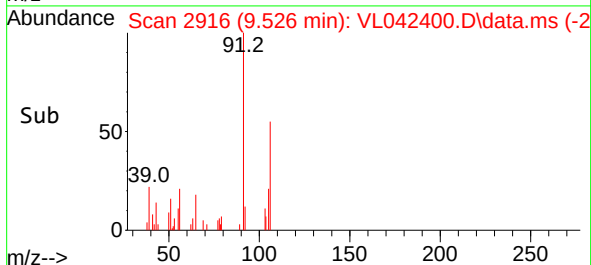
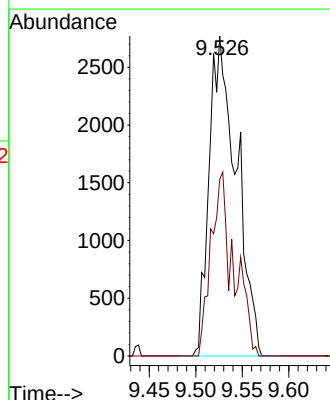
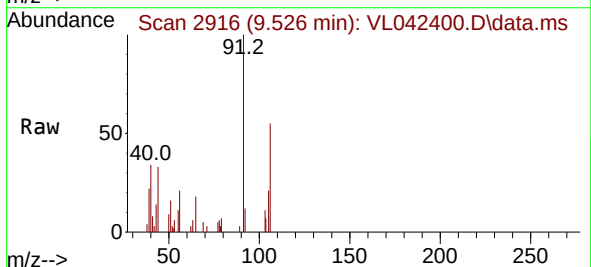
RT: 9.526 min Scan# 2916

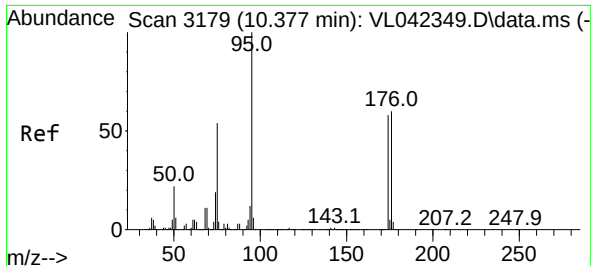
Delta R.T. -0.026 min

Lab File: VL042400.D

Acq: 21 Apr 2025 12:07

Tgt Ion	Ratio	Lower	Upper
91	100		
106	0.0	35.4	53.2





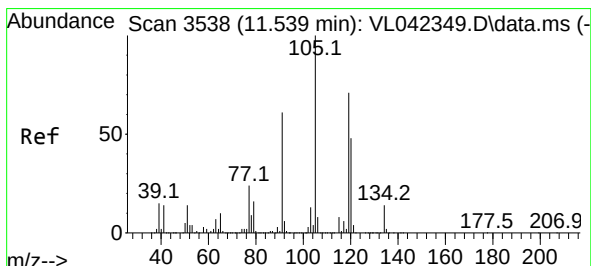
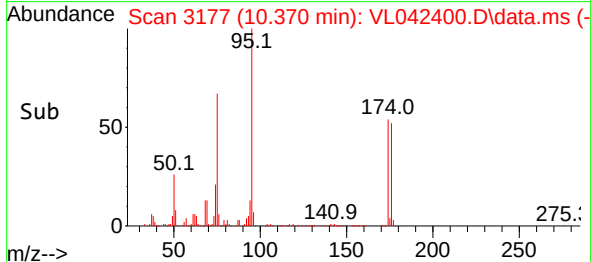
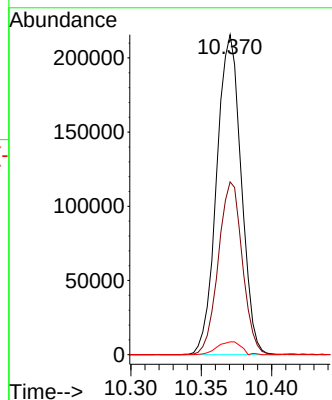
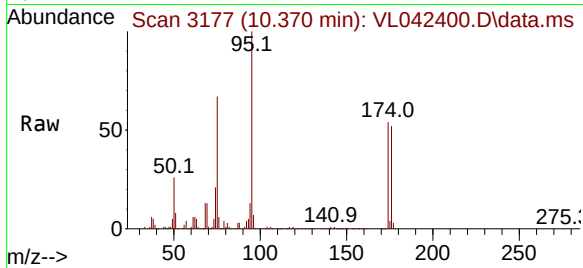
#68
 1-Bromo-4-Fluorobenzene
 Concen: 10.470 ppbv
 RT: 10.370 min Scan# 3179
 Delta R.T. -0.006 min
 Lab File: VL042400.D
 Acq: 21 Apr 2025 12:07

Instrument :
 MSVOA_L
 ClientSampleId :
 SS2R

Tgt Ion: 95 Resp: 261130
 Ion Ratio Lower Upper
 95 100
 174 53.0 48.5 72.7
 175 4.0 3.7 5.5

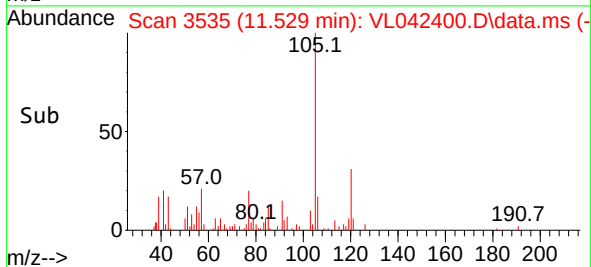
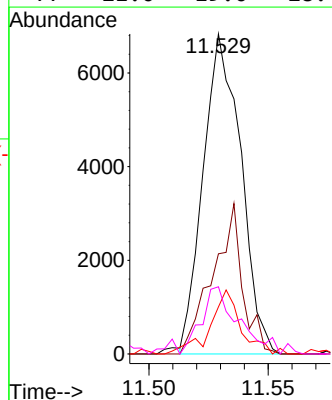
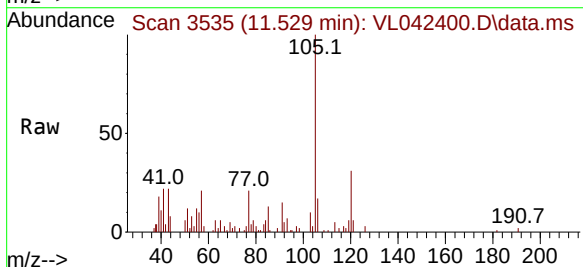
Manual Integrations
 APPROVED

Reviewed By :Semsettin Yesilyurt 04/22/2025
 Supervised By :Mahesh Dadoda 04/22/2025



#74
 1,2,4-Trimethylbenzene
 Concen: 0.122 ppbv
 RT: 11.529 min Scan# 3535
 Delta R.T. -0.010 min
 Lab File: VL042400.D
 Acq: 21 Apr 2025 12:07

Tgt Ion:105 Resp: 7515
 Ion Ratio Lower Upper
 105 100
 120 31.3 38.1 57.1#
 91 12.9 48.8 73.2#
 77 21.0 19.0 28.4



Report of Analysis

Client:	GFE LLC	Date Collected:	04/12/25
Project:	416 Clinton St Hempstead	Date Received:	04/14/25
Client Sample ID:	SS3	SDG No.:	Q1801
Lab Sample ID:	Q1801-02	Matrix:	Air
Analytical Method:	TO-15	Test:	TO-15
Sample Wt/Vol:	400	Units:	mL

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
VL042401.D	4		04/21/25 12:51	VL042125

CAS Number	Parameter	Conc. ppbv	Conc. ug/M3	Qualifier	MDL	LOQ / CRQL	Units
TARGETS							
75-71-8	Dichlorodifluoromethane	0.84	4.15	U	4.15	9.89	ug/m3
74-87-3	Chloromethane	0.39	0.81	U	0.81	4.13	ug/m3
75-01-4	Vinyl Chloride	0.060	0.15	U	0.15	0.31	ug/m3
74-83-9	Bromomethane	0.56	2.17	U	2.17	7.77	ug/m3
75-00-3	Chloroethane	0.68	1.79	U	1.79	5.28	ug/m3
109-99-9	Tetrahydrofuran	0.52	1.53	U	1.53	5.90	ug/m3
75-69-4	Trichlorofluoromethane	0.64	3.60	U	3.60	11.2	ug/m3
76-13-1	1,1,2-Trichlorotrifluoroethane	0.64	4.91	U	4.91	15.3	ug/m3
76-14-2	Dichlorotetrafluoroethane	0.35	2.45	U	2.45	14.0	ug/m3
75-65-0	tert-Butyl alcohol	0.76	2.30	U	2.30	6.06	ug/m3
142-82-5	Heptane	0.44	1.80	U	1.80	8.20	ug/m3
75-35-4	1,1-Dichloroethene	0.56	2.22	U	2.22	7.93	ug/m3
67-64-1	Acetone	5.90	14.0		2.28	4.75	ug/m3
75-15-0	Carbon Disulfide	0.64	1.99	U	1.99	6.23	ug/m3
1634-04-4	Methyl tert-Butyl Ether	0.48	1.73	U	1.73	7.21	ug/m3
75-09-2	Methylene Chloride	0.76	2.64	U	2.64	6.95	ug/m3
156-60-5	trans-1,2-Dichloroethene	0.76	3.01	U	3.01	7.93	ug/m3
75-34-3	1,1-Dichloroethane	0.52	2.10	U	2.10	8.09	ug/m3
110-82-7	Cyclohexane	0.88	3.03	U	3.03	6.88	ug/m3
78-93-3	2-Butanone	0.52	1.53	J	1.42	5.90	ug/m3
56-23-5	Carbon Tetrachloride	0.040	0.25	U	0.25	0.75	ug/m3
156-59-2	cis-1,2-Dichloroethene	0.36	1.43	U	1.43	7.93	ug/m3
67-66-3	Chloroform	0.50	2.44	J	1.61	9.77	ug/m3
71-55-6	1,1,1-Trichloroethane	0.040	0.22	U	0.22	0.65	ug/m3
540-84-1	2,2,4-Trimethylpentane	0.40	1.87	U	1.87	9.34	ug/m3
71-43-2	Benzene	0.35	1.12	U	1.12	6.39	ug/m3
107-06-2	1,2-Dichloroethane	0.36	1.46	U	1.46	8.09	ug/m3
79-01-6	Trichloroethene	0.070	0.38	U	0.38	0.64	ug/m3
78-87-5	1,2-Dichloropropane	0.44	2.03	U	2.03	9.24	ug/m3
75-27-4	Bromodichloromethane	0.31	2.08	U	2.08	13.4	ug/m3
108-10-1	4-Methyl-2-Pentanone	0.37	1.52	U	1.52	8.20	ug/m3
108-88-3	Toluene	1.40	5.28	J	1.66	7.54	ug/m3
10061-02-6	t-1,3-Dichloropropene	0.24	1.09	U	1.09	9.08	ug/m3
10061-01-5	cis-1,3-Dichloropropene	0.26	1.18	U	1.18	9.08	ug/m3
79-00-5	1,1,2-Trichloroethane	0.28	1.53	U	1.53	10.9	ug/m3

Report of Analysis

Client:	GFE LLC	Date Collected:	04/12/25
Project:	416 Clinton St Hempstead	Date Received:	04/14/25
Client Sample ID:	SS3	SDG No.:	Q1801
Lab Sample ID:	Q1801-02	Matrix:	Air
Analytical Method:	TO-15	Test:	TO-15
Sample Wt/Vol:	400	Units:	mL

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
VL042401.D	4		04/21/25 12:51	VL042125

CAS Number	Parameter	Conc. ppbv	Conc. ug/M3	Qualifier	MDL	LOQ / CRQL	Units
124-48-1	Dibromochloromethane	0.26	2.22	U	2.22	17.0	ug/m3
106-93-4	1,2-Dibromoethane	0.29	2.23	U	2.23	3.07	ug/m3
127-18-4	Tetrachloroethene	4.20	28.5		0.41	0.81	ug/m3
108-90-7	Chlorobenzene	0.30	1.38	U	1.38	9.21	ug/m3
100-41-4	Ethyl Benzene	0.48	2.08	U	2.08	8.69	ug/m3
179601-23-1	m/p-Xylene	1.00	4.34	J	3.65	17.4	ug/m3
95-47-6	o-Xylene	0.48	2.08	U	2.08	8.69	ug/m3
100-42-5	Styrene	0.44	1.87	U	1.87	8.52	ug/m3
75-25-2	Bromoform	0.23	2.38	U	2.38	20.7	ug/m3
79-34-5	1,1,2,2-Tetrachloroethane	0.040	0.27	U	0.27	0.82	ug/m3
95-49-8	2-Chlorotoluene	0.44	2.28	U	2.28	10.4	ug/m3
108-67-8	1,3,5-Trimethylbenzene	0.44	2.16	U	2.16	9.83	ug/m3
95-63-6	1,2,4-Trimethylbenzene	0.90	4.42	J	1.52	9.83	ug/m3
541-73-1	1,3-Dichlorobenzene	0.32	1.92	U	1.92	12.0	ug/m3
106-46-7	1,4-Dichlorobenzene	0.21	1.26	U	1.26	12.0	ug/m3
95-50-1	1,2-Dichlorobenzene	0.30	1.80	U	1.80	12.0	ug/m3
120-82-1	1,2,4-Trichlorobenzene	0.34	2.52	U	2.52	14.8	ug/m3
87-68-3	Hexachloro-1,3-Butadiene	0.34	3.63	U	3.63	21.3	ug/m3
106-99-0	1,3-Butadiene	0.52	1.15	U	1.15	4.42	ug/m3
91-20-3	Naphthalene	0.30	1.57	U	1.57	2.10	ug/m3
622-96-8	4-Ethyltoluene	0.48	2.36	U	2.36	9.83	ug/m3
110-54-3	Hexane	4.10	14.4		1.55	7.05	ug/m3
107-05-1	Allyl Chloride	0.60	1.88	U	1.88	6.26	ug/m3
123-91-1	1,4-Dioxane	0.84	3.03	U	3.03	7.21	ug/m3
80-62-6	Methyl Methacrylate	0.34	1.39	U	1.39	8.19	ug/m3
SURROGATES							
460-00-4	1-Bromo-4-Fluorobenzene	10.6			65 - 135	106%	SPK: 10
INTERNAL STANDARDS							
74-97-5	Bromochloromethane	122000		2.787			
540-36-3	1,4-Difluorobenzene	338000		3.958			
3114-55-4	Chlorobenzene-d5	287000		8.878			

Report of Analysis

Client:	GFE LLC	Date Collected:	04/12/25
Project:	416 Clinton St Hempstead	Date Received:	04/14/25
Client Sample ID:	SS3	SDG No.:	Q1801
Lab Sample ID:	Q1801-02	Matrix:	Air
Analytical Method:	TO-15	Test:	TO-15
Sample Wt/Vol:	400	Units:	mL

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
VL042401.D	4		04/21/25 12:51	VL042125

CAS Number	Parameter	Conc. ppbv	Conc. ug/M3	Qualifier	MDL	LOQ / CRQL	Units
------------	-----------	---------------	----------------	-----------	-----	------------	-------

U = Not Detected
 RL = Reporting Limit
 MDL = Method Detection Limit
 E = Value Exceeds Calibration Range
 D = Dilution

J = Estimated Value
 B = Analyte Found in Associated Method Blank
 N = Presumptive Evidence of a Compound
 * = Values outside of QC limits
 Q = indicates LCS control criteria did not meet requirements

Data Path : Z:\voasrv\HPCHEM1\MSVOA_L\Data\VL042125\
 Data File : VL042401.D
 Acq On : 21 Apr 2025 12:51
 Operator : SY/MD
 Sample : Q1801-02 4X
 Misc : 400mL/MSVOA_L
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_L
 ClientSampleId :
 SS3

Manual Integrations
 APPROVED

Reviewed By :Semsettin Yesilyurt 04/22/2025
 Supervised By :Mahesh Dadoda 04/22/2025

Quant Time: Apr 22 02:53:05 2025

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_L\methods\VL041725AIR.M

Quant Title : AIR ANALYSIS BY METHOD TO-15 Instrument: MSVOA_LFri Aug 2

QLast Update : Fri Apr 18 02:14:00 2025

Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) Bromochloromethane	2.787	49	122041	10.000	ppbv	0.00
33) 1,4-Difluorobenzene	3.958	114	337832	10.000	ppbv	0.00
55) Chlorobenzene-d5	8.878	117	286610	10.000	ppbv	0.00
System Monitoring Compounds						
68) 1-Bromo-4-Fluorobenzene	10.374	95	248812	10.564	ppbv	0.00
Spiked Amount	10.000	Range	65 - 135	Recovery	=	105.600%
Target Compounds						
					Qvalue	
2) Dichlorodifluoromethane	1.499	85	2467	0.114	ppbv	97
9) Propene	1.486	41	1848	0.198	ppbv	88
13) Ethanol	1.722	45	11547	7.769	ppbv #	24
15) Acetone	1.835	43	29636	1.483	ppbv	97
17) tert-Butyl alcohol	2.042	59	3820	0.154	ppbv #	71
19) Isopropyl Alcohol	1.887	45	9166	0.826	ppbv #	17
20) Methylene Chloride	2.059	84	1438	0.184	ppbv #	86
23) Vinyl Acetate	2.463	43	6182	0.764	ppbv #	74
26) Hexane	2.822	57	20551	1.031	ppbv #	24
29) Chloroform	2.848	83	3530	0.125	ppbv #	67
34) 2-Butanone	2.563	43	3442	0.129	ppbv #	77
52) Tetrachloroethene	8.221	164	14958	1.062	ppbv	89
53) Toluene	6.933	91	15319m	0.350	ppbv	
59) m/p-Xylene	9.522	91	11844m	0.259	ppbv	
60) o-Xylene	9.962	91	5177	0.112	ppbv	95
74) 1,2,4-Trimethylbenzene	11.535	105	13013	0.224	ppbv #	63

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

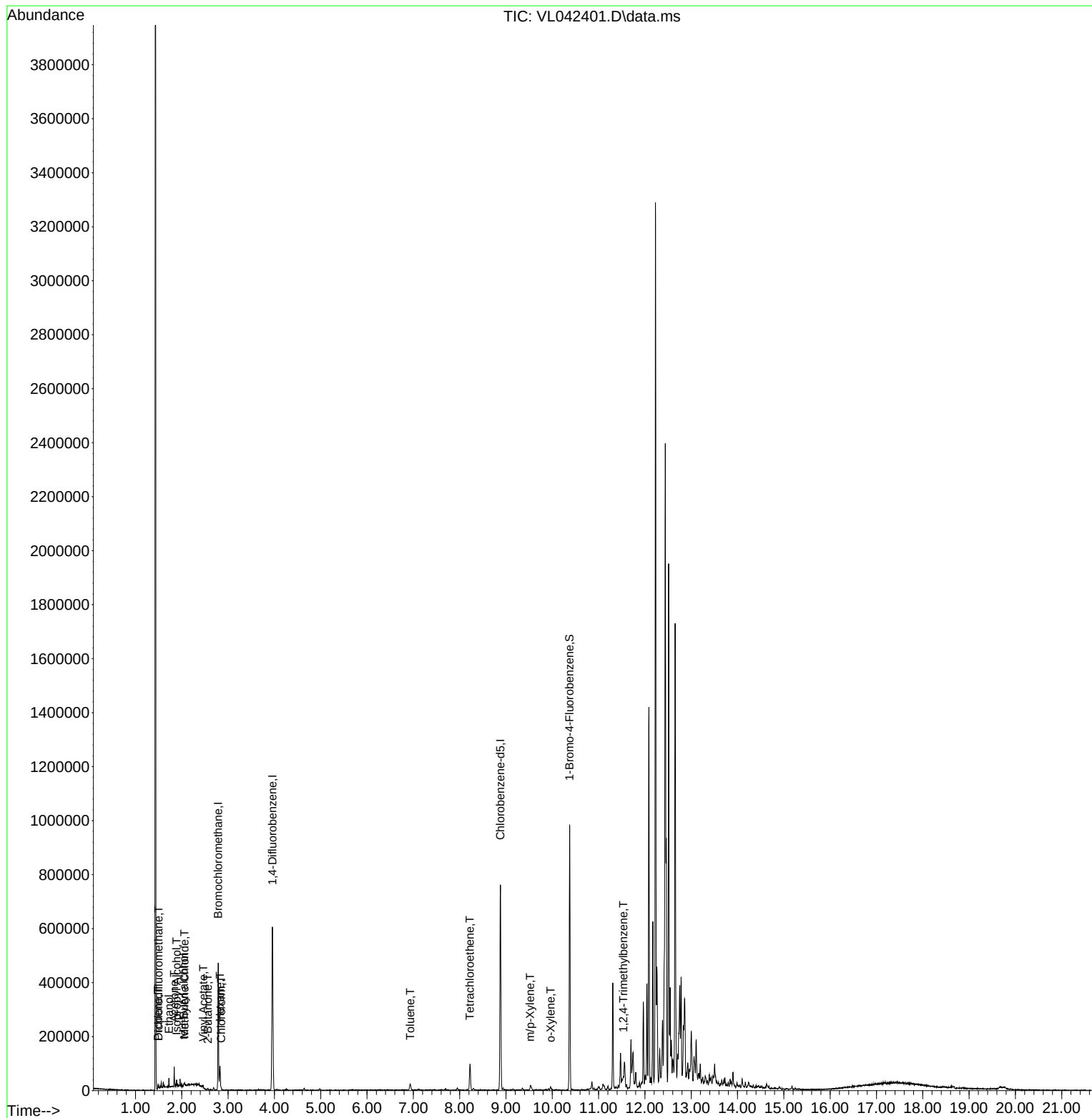
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Data File : VL042401.D
Acq On : 21 Apr 2025 12:51
Operator : SY/MD
Sample : Q1801-02 4X
Misc : 400mL/MSVOA_L
ALS Vial : 1 Sample Multiplier: 1

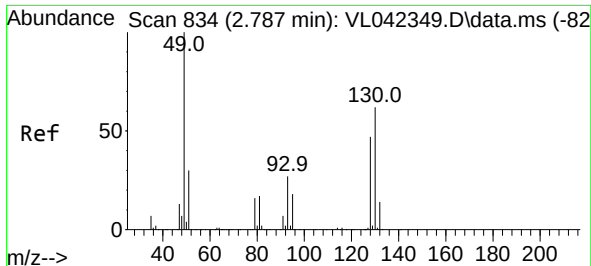
Instrument :
MSVOA_L
ClientSampleId :
SS3

Quant Time: Apr 22 02:53:05 2025
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_L\methods\VL041725AIR.M
Quant Title : AIR ANALYSIS BY METHOD TO-15 Instrument: MSVOA_L Fri Aug
QLast Update : Fri Apr 18 02:14:00 2025
Response via : Initial Calibration

Manual Integrations
APPROVED

Reviewed By :Semsettin Yesilyurt 04/22/2025
Supervised By :Mahesh Dadoda 04/22/2025





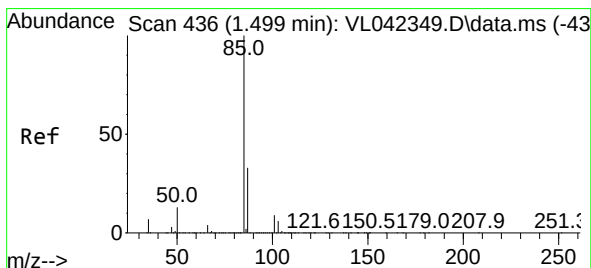
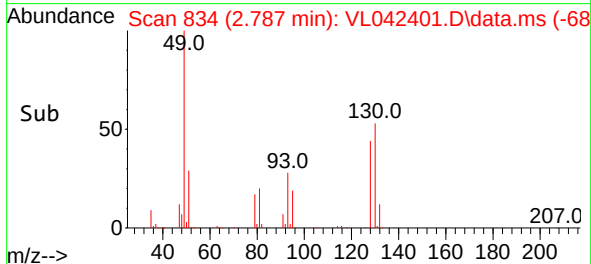
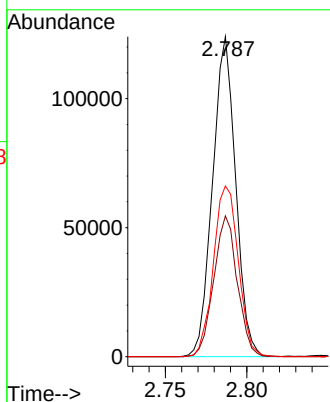
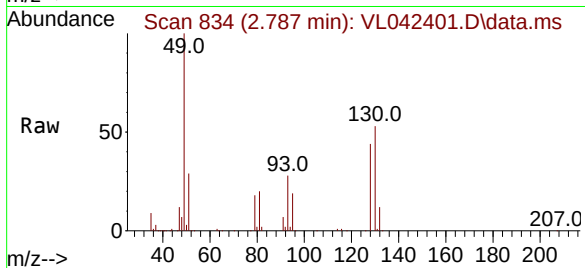
#1
Bromochloromethane
Concen: 10.000 ppbv
RT: 2.787 min Scan# 81
Delta R.T. -0.000 min
Lab File: VL042401.D
Acq: 21 Apr 2025 12:51

Instrument :
MSVOA_L
ClientSampleId :
SS3

Tgt Ion: 49 Resp: 12204
Ion Ratio Lower Upper
49 100
128 44.9 22.8 68.4
130 57.5 29.3 88.0

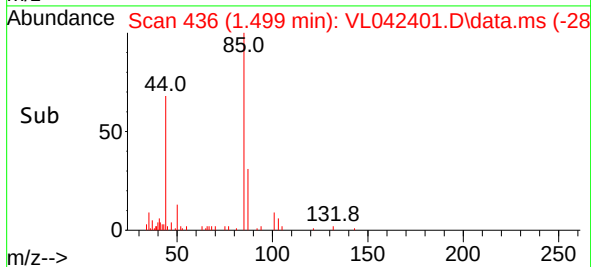
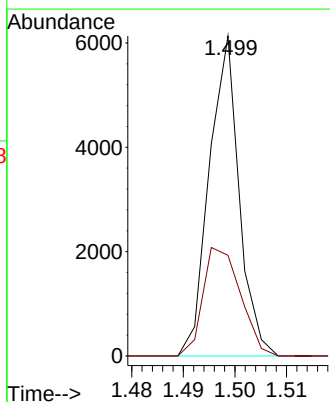
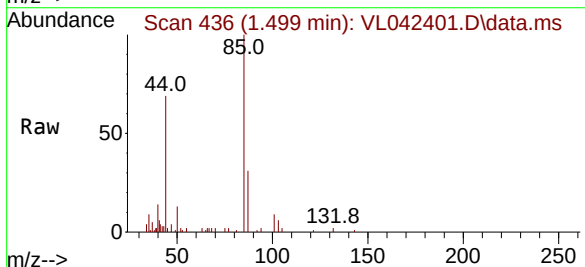
Manual Integrations
APPROVED

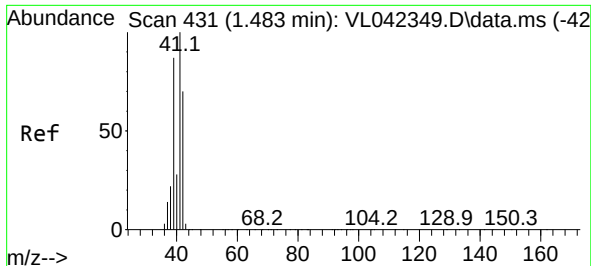
Reviewed By :Semsettin Yesilyurt 04/22/2025
Supervised By :Mahesh Dadoda 04/22/2025



#2
Dichlorodifluoromethane
Concen: 0.114 ppbv
RT: 1.499 min Scan# 436
Delta R.T. -0.000 min
Lab File: VL042401.D
Acq: 21 Apr 2025 12:51

Tgt Ion: 85 Resp: 2467
Ion Ratio Lower Upper
85 100
87 31.5 26.6 39.8





#9

Propene

Concen: 0.198 ppbv

RT: 1.486 min Scan# 41

Delta R.T. 0.003 min

Lab File: VL042401.D

Acq: 21 Apr 2025 12:51

Instrument :

MSVOA_L

ClientSampleId :

SS3

Tgt Ion: 41 Resp: 184

Ion Ratio Lower Upper

41 100

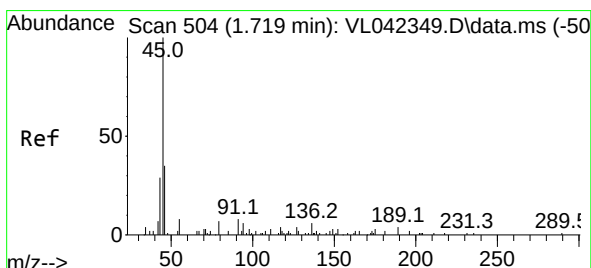
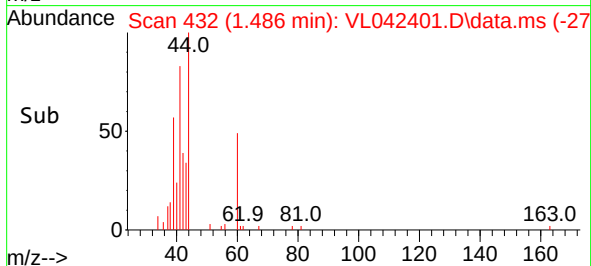
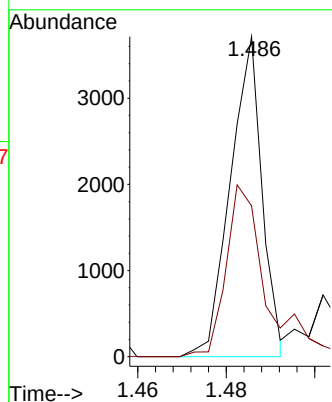
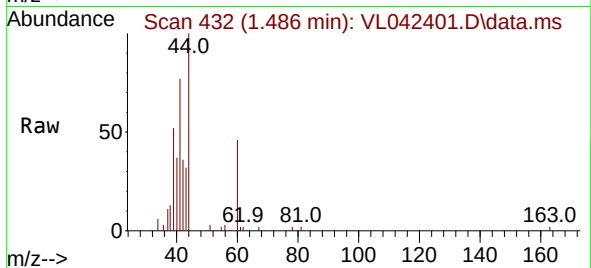
42 80.0 56.2 84.4

Manual Integrations

APPROVED

Reviewed By :Semsettin Yesilyurt 04/22/2025

Supervised By :Mahesh Dadoda 04/22/2025



#13

Ethanol

Concen: 7.769 ppbv

RT: 1.722 min Scan# 505

Delta R.T. 0.003 min

Lab File: VL042401.D

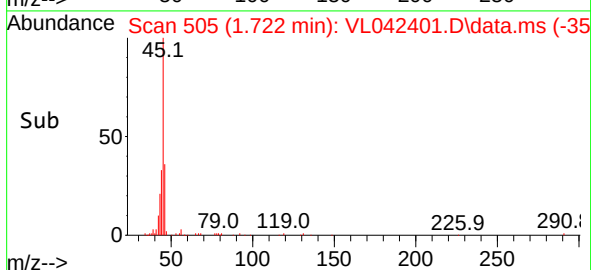
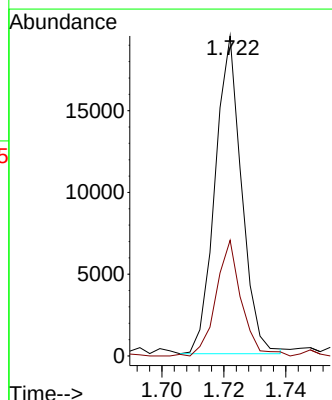
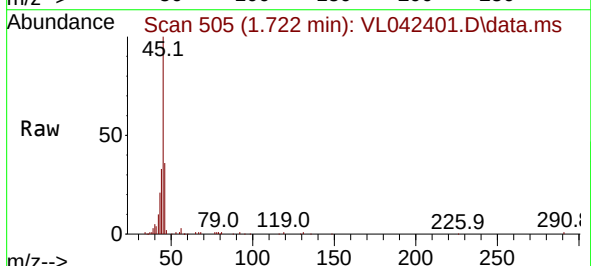
Acq: 21 Apr 2025 12:51

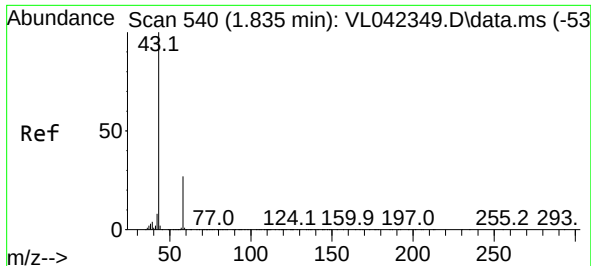
Tgt Ion: 45 Resp: 11547

Ion Ratio Lower Upper

45 100

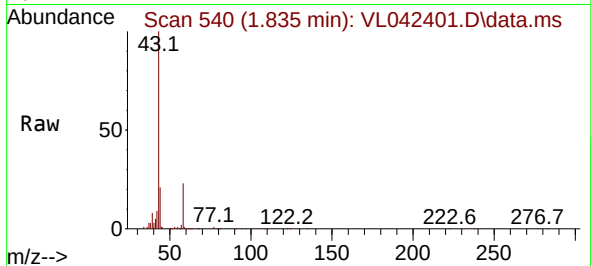
46 0.0 21.9 87.7#





#15
Acetone
Concen: 1.483 ppbv
RT: 1.835 min Scan# 540
Delta R.T. -0.000 min
Lab File: VL042401.D
Acq: 21 Apr 2025 12:51

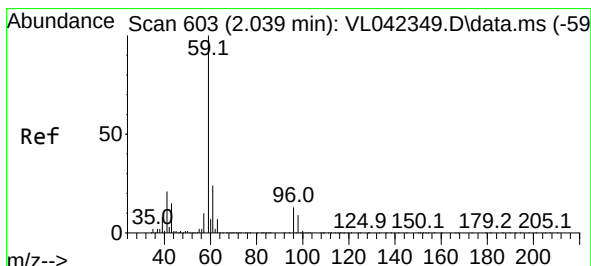
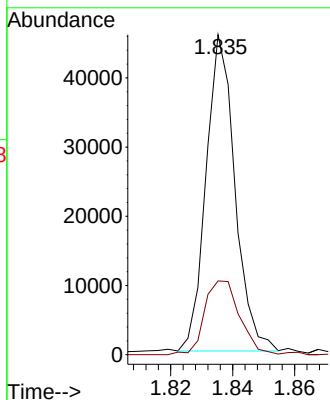
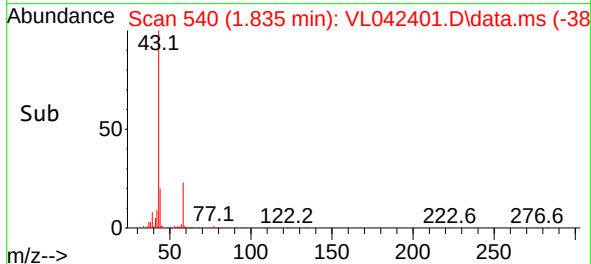
Instrument :
MSVOA_L
ClientSampleId :
SS3



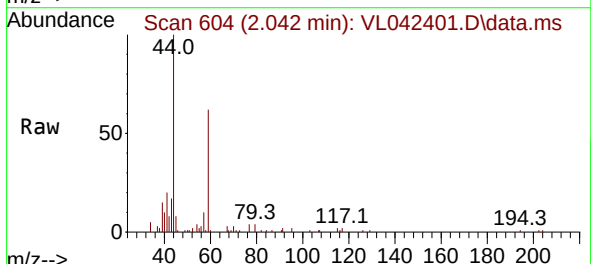
Tgt Ion: 43 Resp: 29630
Ion Ratio Lower Upper
43 100
58 29.8 22.4 33.6

Manual Integrations
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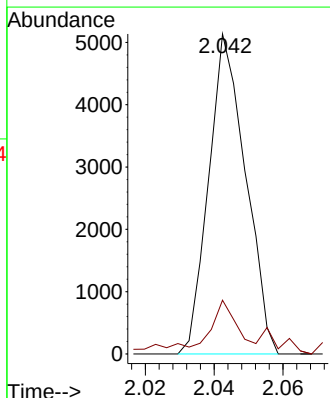
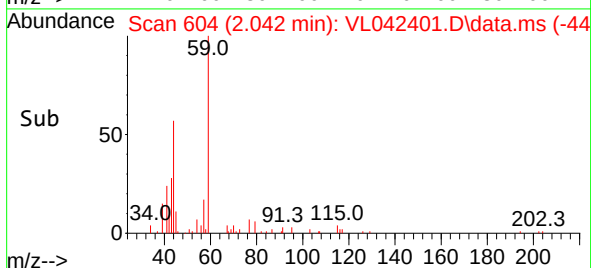
Reviewed By :Semsettin Yesilyurt 04/22/2025
Supervised By :Mahesh Dadoda 04/22/2025

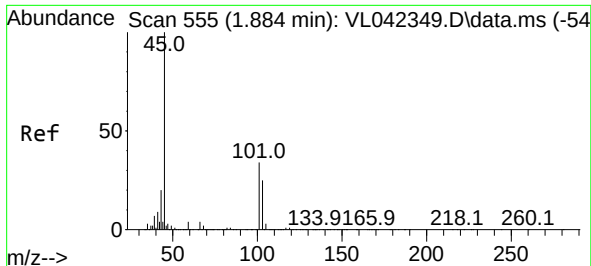


#17
tert-Butyl alcohol
Concen: 0.154 ppbv
RT: 2.042 min Scan# 604
Delta R.T. 0.003 min
Lab File: VL042401.D
Acq: 21 Apr 2025 12:51



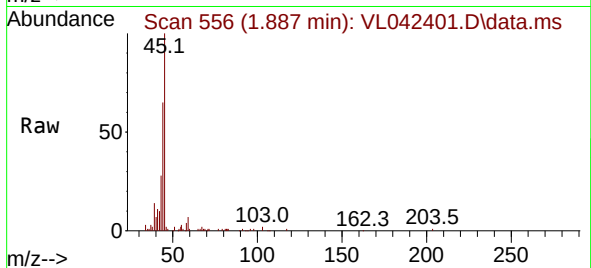
Tgt Ion: 59 Resp: 3820
Ion Ratio Lower Upper
59 100
57 22.0 8.7 13.1#





#19
Isopropyl Alcohol
Concen: 0.826 ppbv
RT: 1.887 min Scan# 516
Delta R.T. 0.003 min
Lab File: VL042401.D
Acq: 21 Apr 2025 12:51

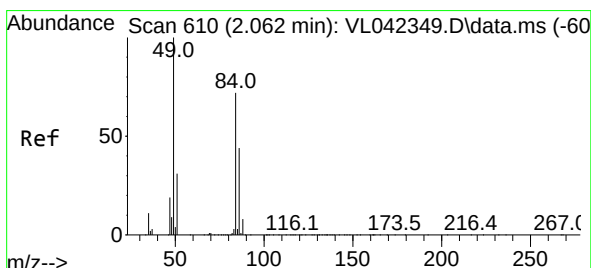
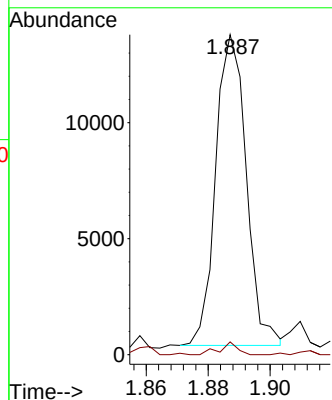
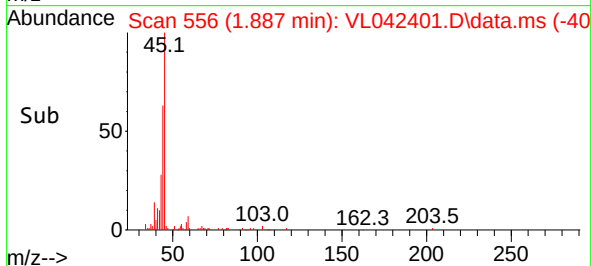
Instrument :
MSVOA_L
ClientSampleId :
SS3



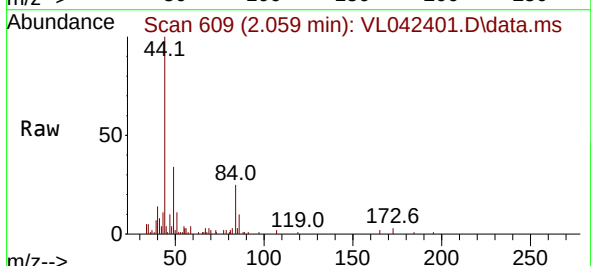
Tgt Ion: 45 Resp: 9160
Ion Ratio Lower Upper
45 100
58 96.2 34.5 51.7

Manual Integrations
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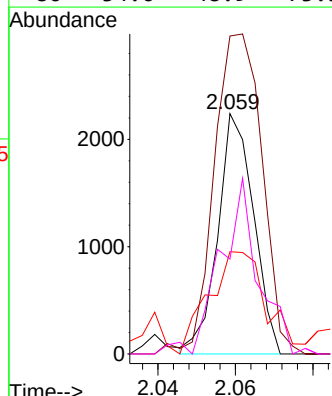
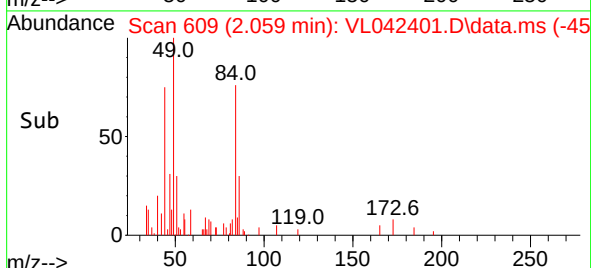
Reviewed By :Semsettin Yesilyurt 04/22/2025
Supervised By :Mahesh Dadoda 04/22/2025

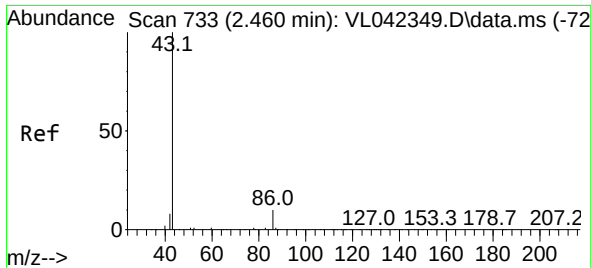


#20
Methylene Chloride
Concen: 0.184 ppbv
RT: 2.059 min Scan# 609
Delta R.T. -0.003 min
Lab File: VL042401.D
Acq: 21 Apr 2025 12:51



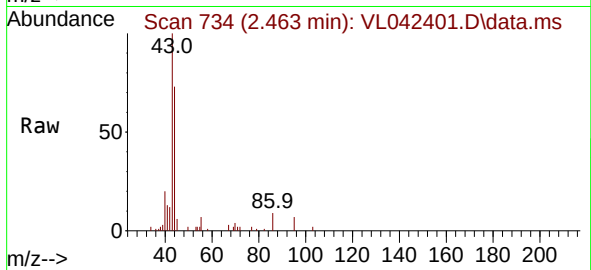
Tgt Ion: 84 Resp: 1438
Ion Ratio Lower Upper
84 100
49 130.0 111.6 167.4
51 42.5 35.1 52.7
86 34.6 48.9 73.3





#23
 Vinyl Acetate
 Concen: 0.764 ppbv
 RT: 2.463 min Scan# 71
 Delta R.T. 0.003 min
 Lab File: VL042401.D
 Acq: 21 Apr 2025 12:51

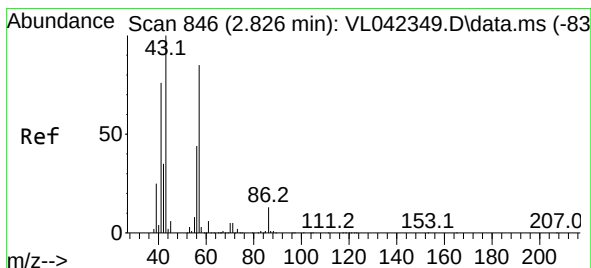
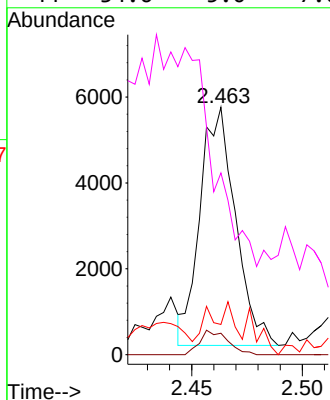
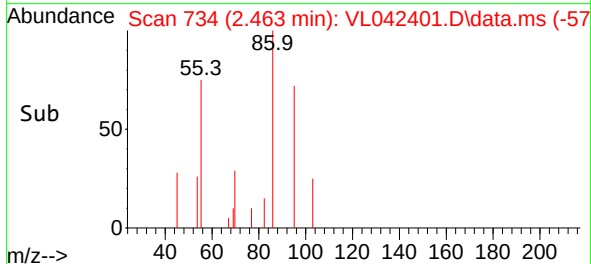
Instrument :
 MSVOA_L
 ClientSampleId :
 SS3



Tgt Ion: 43 Resp: 618
 Ion Ratio Lower Upper
 43 100
 86 9.0 6.5 9.7
 42 12.7 7.8 11.8
 44 34.6 5.0 7.6#

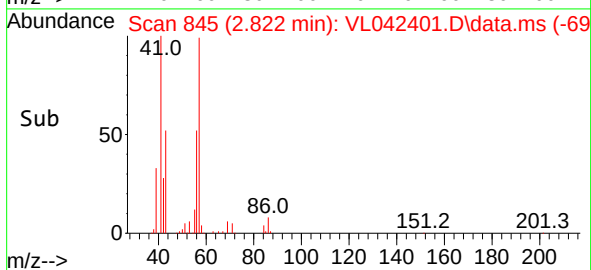
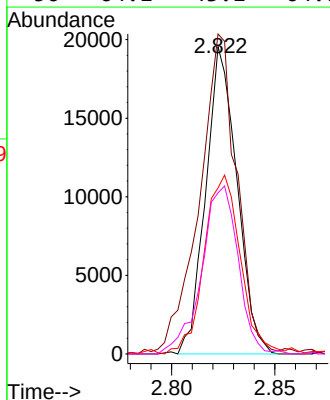
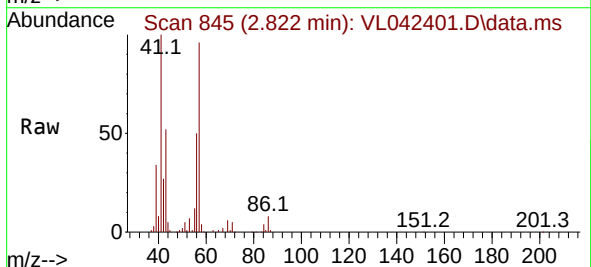
Manual Integrations
 APPROVED

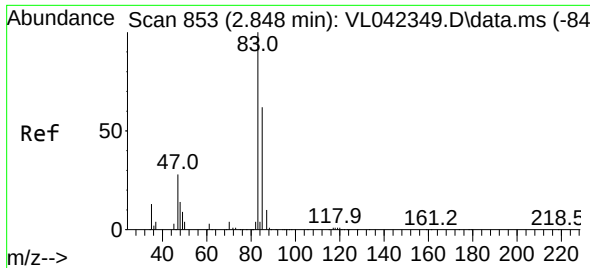
Reviewed By :Semsettin Yesilyurt 04/22/2025
 Supervised By :Mahesh Dadoda 04/22/2025



#26
 Hexane
 Concen: 1.031 ppbv
 RT: 2.822 min Scan# 845
 Delta R.T. -0.003 min
 Lab File: VL042401.D
 Acq: 21 Apr 2025 12:51

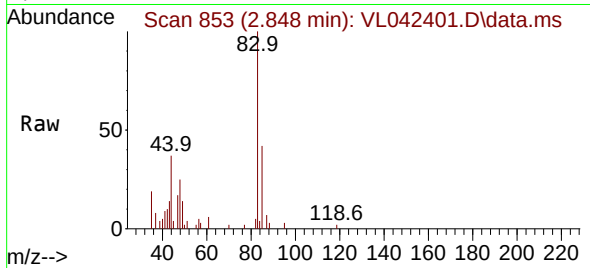
Tgt Ion: 57 Resp: 20551
 Ion Ratio Lower Upper
 57 100
 41 126.8 77.9 116.9#
 43 70.4 208.4 312.6#
 56 64.1 43.1 64.7





#29
Chloroform
Concen: 0.125 ppbv
RT: 2.848 min Scan# 853
Delta R.T. -0.000 min
Lab File: VL042401.D
Acq: 21 Apr 2025 12:51

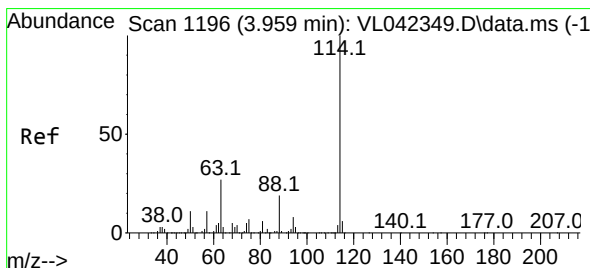
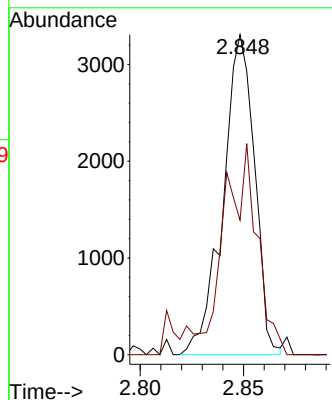
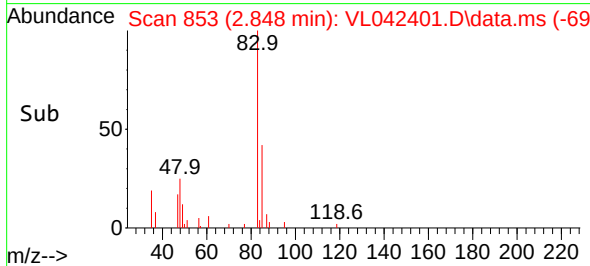
Instrument :
MSVOA_L
ClientSampleId :
SS3



Tgt Ion: 83 Resp: 3530
Ion Ratio Lower Upper
83 100
85 37.3 50.0 75.0

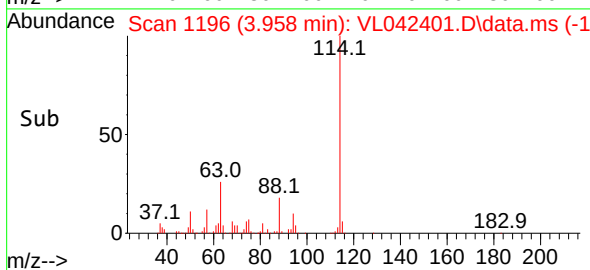
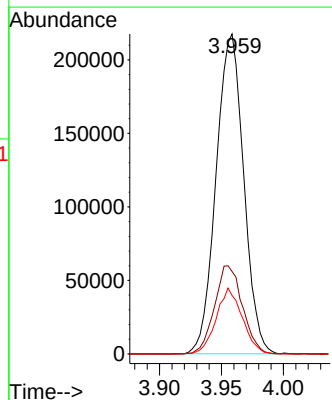
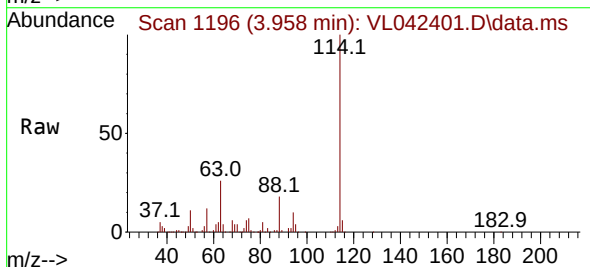
Manual Integrations
APPROVED

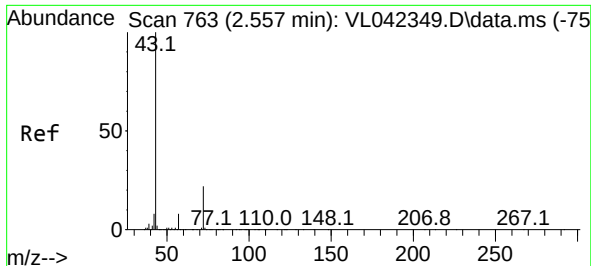
Reviewed By :Semsettin Yesilyurt 04/22/2025
Supervised By :Mahesh Dadoda 04/22/2025



#33
1,4-Difluorobenzene
Concen: 10.000 ppbv
RT: 3.958 min Scan# 1196
Delta R.T. -0.000 min
Lab File: VL042401.D
Acq: 21 Apr 2025 12:51

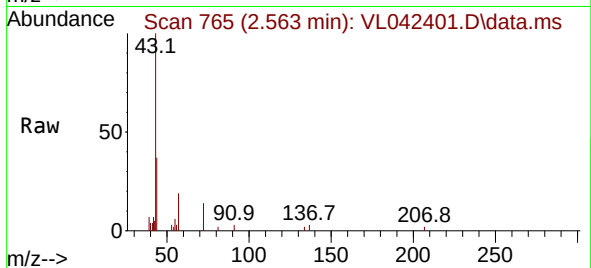
Tgt Ion:114 Resp: 337832
Ion Ratio Lower Upper
114 100
63 28.0 21.8 32.6
88 19.5 15.5 23.3





#34
2-Butanone
Concen: 0.129 ppbv
RT: 2.563 min Scan# 763
Delta R.T. 0.006 min
Lab File: VL042401.D
Acq: 21 Apr 2025 12:51

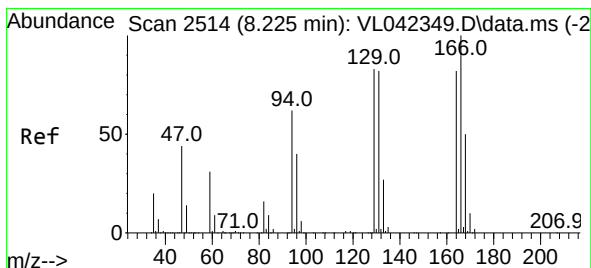
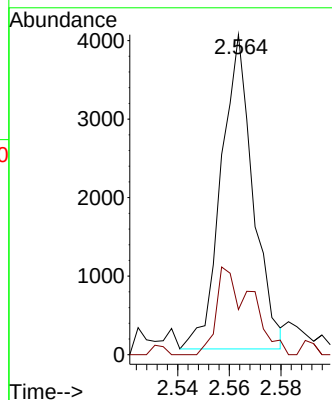
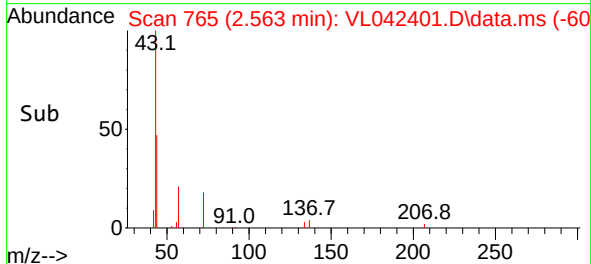
Instrument :
MSVOA_L
ClientSampleId :
SS3



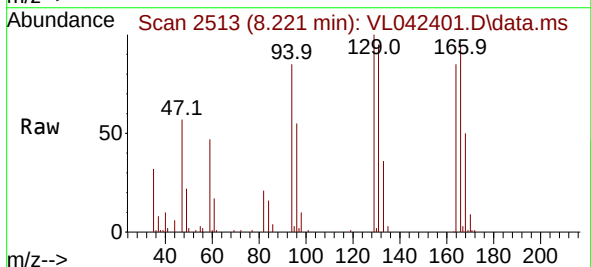
Tgt Ion: 43 Resp: 3442
Ion Ratio Lower Upper
43 100
72 33.6 18.2 27.2

Manual Integrations
APPROVED

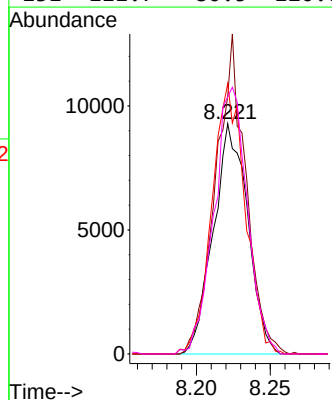
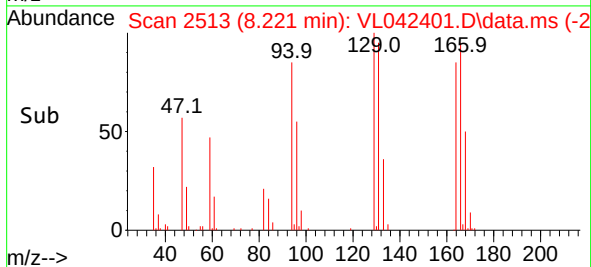
Reviewed By :Semsettin Yesilyurt 04/22/2025
Supervised By :Mahesh Dadoda 04/22/2025

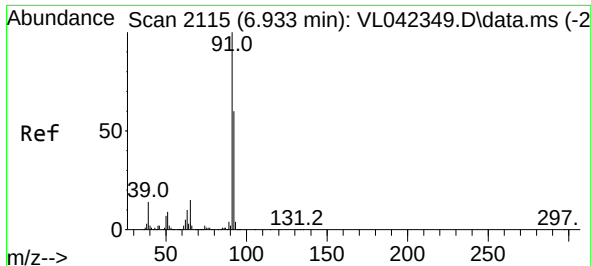


#52
Tetrachloroethene
Concen: 1.062 ppbv
RT: 8.221 min Scan# 2513
Delta R.T. -0.003 min
Lab File: VL042401.D
Acq: 21 Apr 2025 12:51



Tgt Ion:164 Resp: 14958
Ion Ratio Lower Upper
164 100
166 112.0 98.2 147.2
129 116.3 81.8 122.8
131 111.7 80.5 120.7





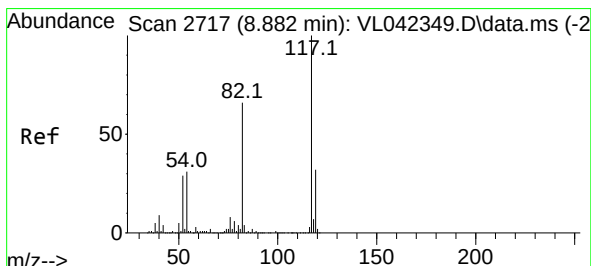
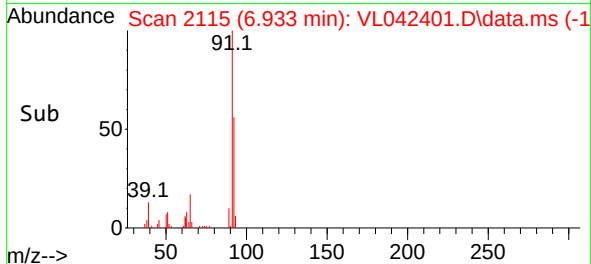
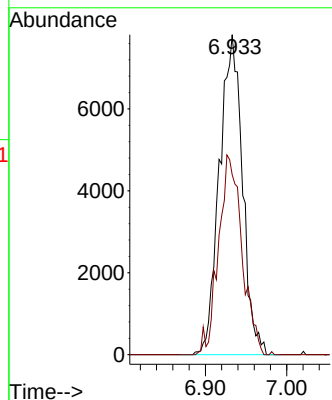
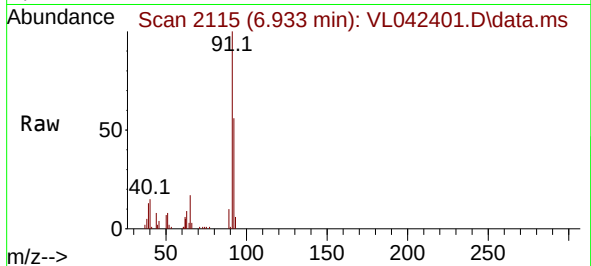
#53
Toluene
Concen: 0.350 ppbv m
RT: 6.933 min Scan# 2115
Delta R.T. -0.000 min
Lab File: VL042401.D
Acq: 21 Apr 2025 12:51

Instrument :
MSVOA_L
ClientSampleId :
SS3

Tgt Ion: 91 Resp: 15319
Ion Ratio Lower Upper
91 100
92 63.6 48.2 72.4

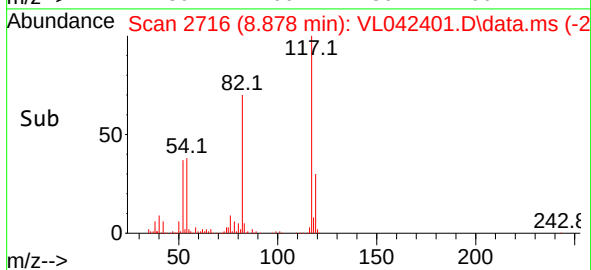
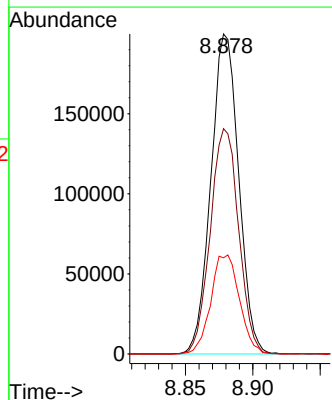
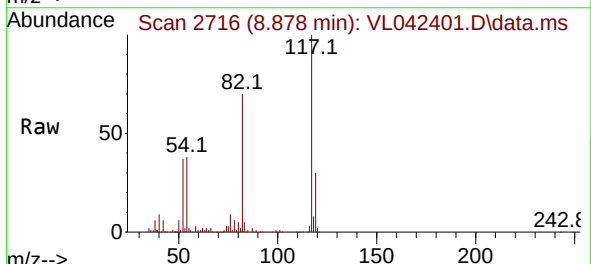
Manual Integrations
APPROVED

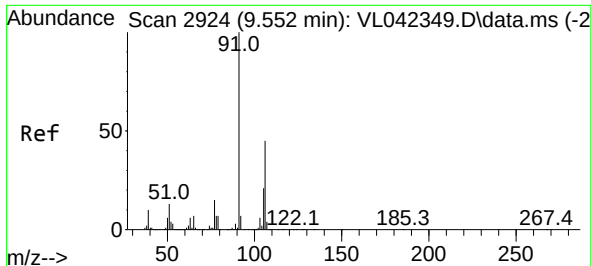
Reviewed By :Semsettin Yesilyurt 04/22/2025
Supervised By :Mahesh Dadoda 04/22/2025



#55
Chlorobenzene-d5
Concen: 10.000 ppbv
RT: 8.878 min Scan# 2716
Delta R.T. -0.003 min
Lab File: VL042401.D
Acq: 21 Apr 2025 12:51

Tgt Ion:117 Resp: 286610
Ion Ratio Lower Upper
117 100
82 71.9 57.8 86.6
119 32.0 25.2 37.8





#59

m/p-Xylene

Concen: 0.259 ppbv m

RT: 9.522 min Scan# 2915

Delta R.T. -0.029 min

Lab File: VL042401.D

Acq: 21 Apr 2025 12:51

Instrument :

MSVOA_L

ClientSampleId :

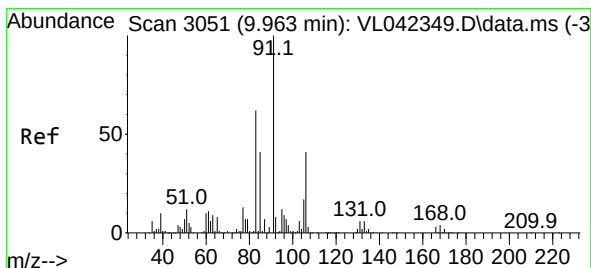
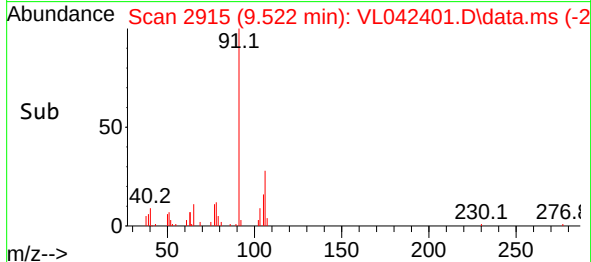
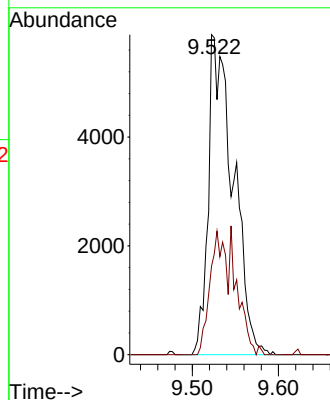
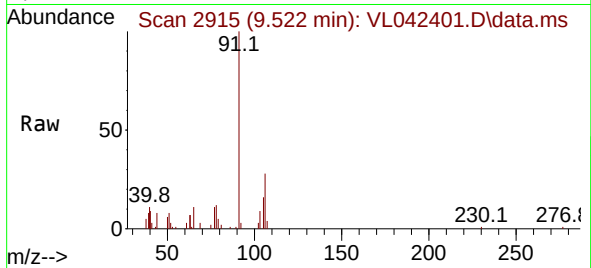
SS3

Tgt Ion: 91 Resp: 1184
 Ion Ratio Lower Upper
 91 100
 106 38.4 35.4 53.2

Manual Integrations

APPROVED

Reviewed By :Semsettin Yesilyurt 04/22/2025
 Supervised By :Mahesh Dadoda 04/22/2025



#60

o-Xylene

Concen: 0.112 ppbv

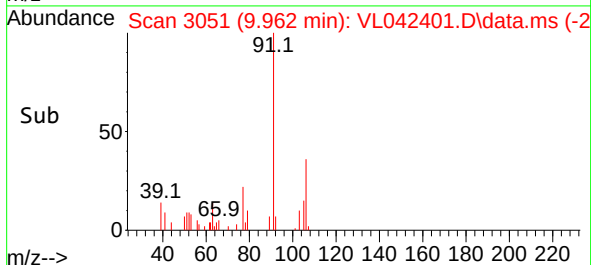
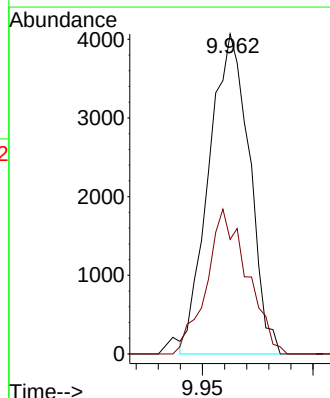
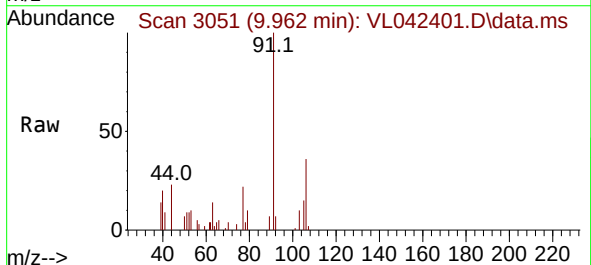
RT: 9.962 min Scan# 3051

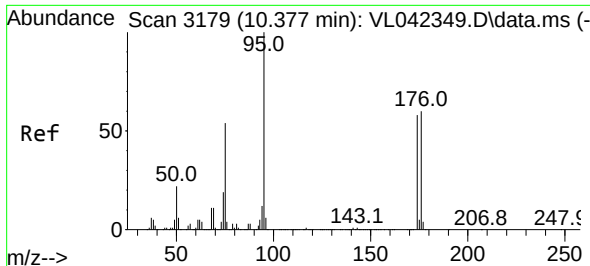
Delta R.T. -0.000 min

Lab File: VL042401.D

Acq: 21 Apr 2025 12:51

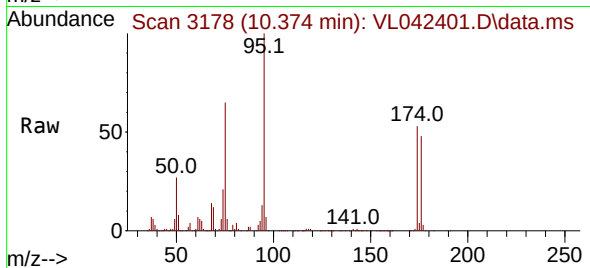
Tgt Ion: 91 Resp: 5177
 Ion Ratio Lower Upper
 91 100
 106 45.5 21.1 63.4





#68
1-Bromo-4-Fluorobenzene
Concen: 10.564 ppbv
RT: 10.374 min Scan# 3179
Delta R.T. -0.003 min
Lab File: VL042401.D
Acq: 21 Apr 2025 12:51

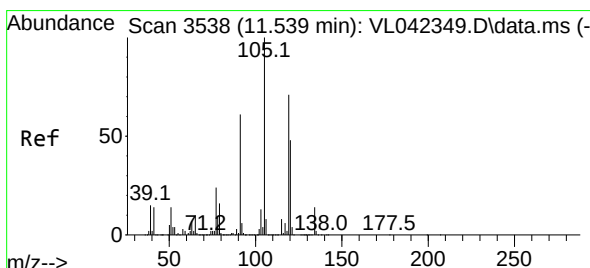
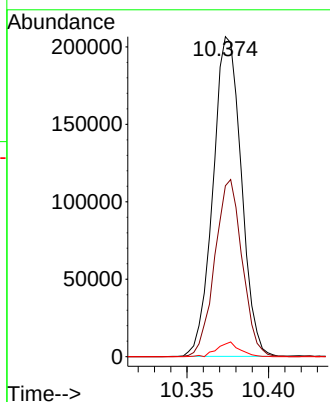
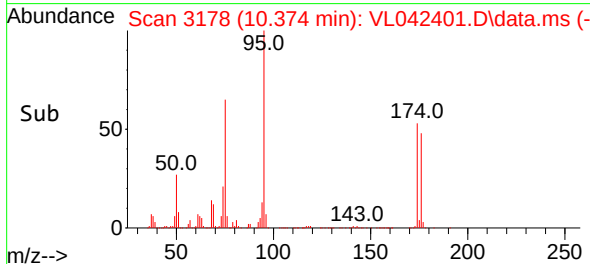
Instrument :
MSVOA_L
ClientSampleId :
SS3



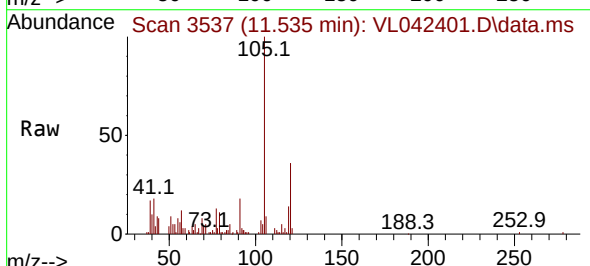
Tgt Ion: 95 Resp: 248812
Ion Ratio Lower Upper
95 100
174 53.8 48.5 72.7
175 3.7 3.7 5.5

Manual Integrations
APPROVED

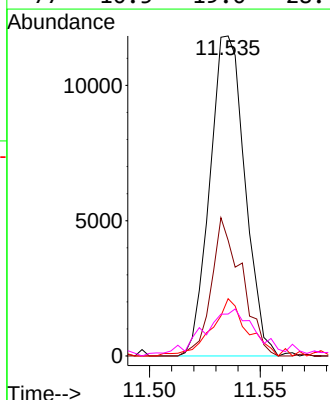
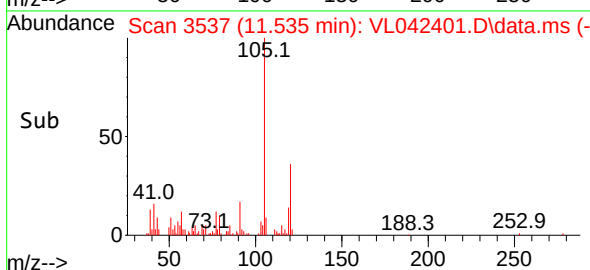
Reviewed By :Semsettin Yesilyurt 04/22/2025
Supervised By :Mahesh Dadoda 04/22/2025



#74
1,2,4-Trimethylbenzene
Concen: 0.224 ppbv
RT: 11.535 min Scan# 3537
Delta R.T. -0.003 min
Lab File: VL042401.D
Acq: 21 Apr 2025 12:51



Tgt Ion:105 Resp: 13013
Ion Ratio Lower Upper
105 100
120 36.1 38.1 57.1#
91 17.9 48.8 73.2#
77 10.5 19.0 28.4#





CALIBRATION SUMMARY

VOLATILE ORGANICS INITIAL CALIBRATION DATA

Lab Name: CHEMTECH Contract: GFEL01
 Lab Code: CHEM Case No.: Q1801 SAS No.: Q1801 SDG No.: Q1801
 Instrument ID: MSVOA_L Calibration Date(s): 04/17/2025 04/17/2025
 Heated Purge: (Y/N) N Calibration Time(s): 09:44 12:54
 GC Column: RTX-1 ID: 0.32 (mm)

LAB FILE ID:		RRF010 = VL042349.D		RRF002 = VL042350.D		RRF001 = VL042351.D		
		RRF0.5 = VL042352.D		RRF0.1 = VL042353.D		RRF.03 = VL042354.D		
COMPOUND	RRF010	RRF002	RRF001	RRF0.5	RRF0.1	RRF.03	RRF	% RSD
Dichlorodifluoromethane	1.658	1.827	1.837	1.831			1.767	5
Chloromethane	0.652	0.701	0.725	0.752			0.699	5.9
Vinyl Chloride	0.638	0.689	0.709	0.731	0.911	1.002	0.761	18.4
Bromomethane	0.304	0.300	0.330	0.397			0.327	12.6
Chloroethane	0.253	0.280	0.287	0.303			0.275	7.8
Tetrahydrofuran	0.409	0.468	0.480	0.518			0.461	9.2
Trichlorofluoromethane	1.567	1.746	1.806	1.684			1.680	6
1,1,2-Trichlorotrifluoroethane	1.281	1.437	1.439	1.389			1.374	5.1
Dichlorotetrafluoroethane	1.432	1.581	1.550	1.556			1.517	4.3
tert-Butyl alcohol	1.804	2.103	2.132	2.263			2.029	9.8
Heptane	1.884	2.209	2.296	2.444			2.153	11.1
1,1-Dichloroethene	0.583	0.637	0.635	0.735			0.642	8.8
Acetone	1.263	1.743	1.895	1.968			1.637	20
Carbon Disulfide	1.668	1.785	1.670	1.903			1.749	5.6
Methyl tert-Butyl Ether	0.940	1.132	1.141	1.134			1.070	8.6
Methylene Chloride	0.498	0.582	0.728	0.875			0.641	24.7
trans-1,2-Dichloroethene	0.664	0.736	0.759	0.814			0.731	8.3
1,1-Dichloroethane	1.284	1.383	1.469	1.452			1.388	5.4
Cyclohexane	1.252	1.347	1.333	1.422			1.320	5.5
2-Butanone	0.687	0.818	0.826	0.886			0.791	9.9
Carbon Tetrachloride	0.770	0.847	0.830	0.832	0.914	1.319	0.905	20.7
cis-1,2-Dichloroethene	1.459	1.593	1.603	1.588			1.553	3.9
Chloroform	2.145	2.392	2.333	2.458			2.309	5.5
1,1,1-Trichloroethane	2.274	2.493	2.549	2.557	2.840	3.110	2.599	11
2,2,4-Trimethylpentane	1.708	1.962	1.916	1.935			1.863	5.8
Benzene	1.002	1.121	1.075	1.146			1.079	5.3
1,2-Dichloroethane	0.592	0.666	0.647	0.691			0.647	5.6
Trichloroethene	0.395	0.447	0.448	0.443	0.505	0.643	0.471	17.6
1,2-Dichloropropane	0.370	0.405	0.413	0.441			0.402	7
Bromodichloromethane	0.852	0.912	0.909	0.943			0.901	3.7

* Compounds with required minimum RRF and maximum %RSD values.
 All other compounds must meet a minimum RRF of 0.010.
 RRF of 1,4-Dioxane = Value should be divide by 1000.

VOLATILE ORGANICS INITIAL CALIBRATION DATA

Lab Name:	<u>CHEMTECH</u>	Contract:	<u>GFEL01</u>
Lab Code:	<u>CHEM</u>	Case No.:	<u>Q1801</u>
Instrument ID:	<u>MSVOA_L</u>	SAS No.:	<u>Q1801</u>
Heated Purge: (Y/N)	<u>N</u>	SDG No.:	<u>Q1801</u>
GC Column:	<u>RTX-1</u>	Calibration Date(s):	<u>04/17/2025</u>
ID:	<u>0.32 (mm)</u>	Calibration Time(s):	<u>09:44</u>
			<u>12:54</u>

LAB FILE ID:		RRF010 = VL042349.D		RRF002 = VL042350.D		RRF001 = VL042351.D		
		RRF0.5 = VL042352.D		RRF0.1 = VL042353.D		RRF.03 = VL042354.D		
COMPOUND	RRF010	RRF002	RRF001	RRF0.5	RRF0.1	RRF.03	RRF	% RSD
4-Methyl-2-Pentanone	1.082	1.435	1.382	1.541			1.314	15.1
Toluene	1.193	1.370	1.295	1.371			1.294	6
t-1,3-Dichloropropene	0.459	0.509	0.510	0.514			0.497	4.6
cis-1,3-Dichloropropene	0.570	0.637	0.617	0.613			0.608	4.1
1,1,2-Trichloroethane	0.376	0.420	0.429	0.423			0.410	5.3
Dibromochloromethane	0.679	0.746	0.721	0.743			0.723	3.7
1,2-Dibromoethane	0.539	0.606	0.605	0.582	0.621		0.586	5.3
Tetrachloroethene	0.357	0.391	0.393	0.420	0.452	0.543	0.417	15.5
Chlorobenzene	0.956	1.068	1.067	1.129			1.040	6.8
Ethyl Benzene	1.832	2.068	2.088	2.029			1.981	5.7
m/p-Xylene	1.456	1.658	1.682	1.685			1.596	6.9
o-Xylene	1.442	1.668	1.697	1.740			1.607	8.3
Styrene	0.726	0.808	0.776	0.792			0.767	4.7
Bromoform	0.616	0.675	0.623	0.622			0.633	3.8
1,1,2,2-Tetrachloroethane	0.873	0.968	0.985	0.990	1.176	1.407	1.042	18.1
2-Chlorotoluene	1.695	1.920	1.924	1.972			1.851	6.7
1,3,5-Trimethylbenzene	1.517	1.841	1.851	1.894			1.735	10.2
1,2,4-Trimethylbenzene	1.671	2.168	2.245	2.348			2.026	15.7
1,3-Dichlorobenzene	0.948	1.119	1.121	1.119			1.053	8.6
1,4-Dichlorobenzene	0.944	1.122	1.115	1.138			1.059	8.7
1,2-Dichlorobenzene	0.904	1.114	1.093	1.085			1.024	9.9
1,2,4-Trichlorobenzene	0.905	1.224	1.117	1.026			1.040	12.8
Hexachloro-1,3-Butadiene	0.863	1.135	1.136	1.163			1.034	14.7
1,3-Butadiene	0.651	0.735	0.836	1.015			0.783	18.9
Naphthalene	2.004	2.909	2.756	2.395	2.433		2.423	15.1
4-Ethyltoluene	1.765	2.054	2.118	2.055			1.965	7.9
1-Bromo-4-Fluorobenzene	0.831	0.822	0.818	0.824	0.814	0.819	0.822	0.7
Hexane	1.424	1.625	1.719	1.899			1.633	11.5
Allyl Chloride	0.973	1.110	1.213	1.276			1.130	10.4
1,4-Dioxane	181.866	232.301	242.757	262.618			220.893	16.3

* Compounds with required minimum RRF and maximum %RSD values.
All other compounds must meet a minimum RRF of 0.010.
RRF of 1,4-Dioxane = Value should be divide by 1000.

VOLATILE ORGANICS INITIAL CALIBRATION DATA

Lab Name: CHEMTECH Contract: GFEL01
 Lab Code: CHEM Case No.: Q1801 SAS No.: Q1801 SDG No.: Q1801
 Instrument ID: MSVOA_L Calibration Date(s): 04/17/2025 04/17/2025
 Heated Purge: (Y/N) N Calibration Time(s): 09:44 12:54
 GC Column: RTX-1 ID: 0.32 (mm)

LAB FILE ID:								
RRF010 = VL042349.D			RRF002 = VL042350.D			RRF001 = VL042351.D		
RRF0.5 = VL042352.D			RRF0.1 = VL042353.D			RRF.03 = VL042354.D		
COMPOUND	RRF010	RRF002	RRF001	RRF0.5	RRF0.1	RRF.03	RRF	% RSD
Methyl Methacrylate	0.430	0.495	0.520	0.545			0.486	10.2

* Compounds with required minimum RRF and maximum %RSD values.
 All other compounds must meet a minimum RRF of 0.010.
 RRF of 1,4-Dioxane = Value should be divide by 1000.

Method Path : Z:\voasrv\HPCHEM1\MSVOA_L\methods\

Method File : VL041725AIR.M

Title : AIR ANALYSIS BY METHOD TO-15 Instrument: MSVOA_L Fri Aug 26 06:05:16 2022

Last Update : Fri Apr 18 02:14:00 2025

Response Via : Initial Calibration

Calibration Files

0.03=VL042354.D 0.1 =VL042353.D 0.5 =VL042352.D 1 =VL042351.D 2 =VL042350.D 10 =VL042349.D 15 =VL042355.D

Compound	0.03	0.1	0.5	1	2	10	15	Avg	%RSD

1) I Bromochloromethane	-----ISTD-----								
2) T Dichlorodifluo...			1.831	1.837	1.827	1.658	1.682	1.767	5.05
3) Chlorodifluoro...			1.876	1.657	1.698	1.573	1.608	1.682	7.01
4) Chloromethane			0.752	0.725	0.701	0.652	0.665	0.699	5.93
5) T Vinyl Chloride	1.002	0.911	0.731	0.709	0.689	0.638	0.647	0.761	18.39
6) T Bromomethane			0.397	0.330	0.300	0.304	0.303	0.327	12.56
7) Chloroethane			0.303	0.287	0.280	0.253	0.255	0.275	7.76
8) T Dichlorotetra...			1.556	1.550	1.581	1.432	1.465	1.517	4.25
9) T Propene			0.880	0.797	0.758	0.690	0.707	0.766	9.95
10) T Heptane			2.444	2.296	2.209	1.884	1.931	2.153	11.14
11) T Trichlorofluor...			1.684	1.806	1.746	1.567	1.597	1.680	5.95
12) T 1,1,2-Trichlor...			1.389	1.439	1.437	1.281	1.323	1.374	5.11
13) Ethanol			0.275	0.139	0.123	0.035	0.036	0.122	80.54
14) T Bromoethene			0.537	0.545	0.530	0.490	0.487	0.518	5.24
15) T Acetone			1.968	1.895	1.743	1.263	1.316	1.637	20.03
16) T 1,3-Butadiene			1.015	0.836	0.735	0.651	0.678	0.783	18.90
17) tert-Butyl alc...			2.263	2.132	2.103	1.804	1.841	2.029	9.76
18) T 1,1-Dichloroet...			0.735	0.635	0.637	0.583	0.618	0.642	8.79
19) T Isopropyl Alcohol			1.050	0.980	0.897	0.821	0.798	0.909	11.66
20) T Methylene Chlo...			0.875	0.728	0.582	0.498	0.523	0.641	24.71
21) T Allyl Chloride			1.276	1.213	1.110	0.973	1.080	1.130	10.44
22) T trans-1,2-Dich...			0.814	0.759	0.736	0.664	0.681	0.731	8.32
23) T Vinyl Acetate			0.656	0.744	0.749	0.555	0.610	0.663	12.72
24) T 1,1-Dichloroet...			1.452	1.469	1.383	1.284	1.352	1.388	5.43
25) T Ethyl Acetate			4.428	4.147	4.118	3.677	3.812	4.036	7.34
26) T Hexane			1.899	1.719	1.625	1.424	1.499	1.633	11.45
27) T Carbon Disulfide			1.903	1.670	1.785	1.668	1.718	1.749	5.63
28) T Methyl tert-Bu...			1.134	1.141	1.132	0.940	1.004	1.070	8.64
29) T Chloroform			2.458	2.333	2.392	2.145	2.215	2.309	5.54
30) T Cyclohexane			1.422	1.333	1.347	1.252	1.249	1.320	5.50
31) T cis-1,2-Dichlo...			1.588	1.603	1.593	1.459	1.522	1.553	3.95
32) T 1,1,1-Trichlor...	3.110	2.840	2.557	2.549	2.493	2.274	2.369	2.599	11.03

33) I 1,4-Difluorobenzene	-----ISTD-----								
34) T 2-Butanone			0.886	0.826	0.818	0.687	0.737	0.791	9.94
35) T Carbon Tetrach...	1.319	0.914	0.832	0.830	0.847	0.770	0.823	0.905	20.70
36) T Benzene			1.146	1.075	1.121	1.002	1.048	1.079	5.31
37) T 1,2-Dichloroet...			0.691	0.647	0.666	0.592	0.641	0.647	5.63
38) T Trichloroethene	0.643	0.505	0.443	0.448	0.447	0.395	0.419	0.471	17.59
39) T 1,2-Dichloropr...			0.441	0.413	0.405	0.370	0.381	0.402	6.97

Method Path : Z:\voasrv\HPCHEM1\MSVOA_L\methods\										
Method File : VL041725AIR.M										
40) T	1,4-Dioxane			0.263	0.243	0.232	0.182	0.185	0.221	16.27
41) T	Tetrahydrofuran			0.518	0.480	0.468	0.409	0.432	0.461	9.20
42) T	Bromodichlorom...			0.943	0.909	0.912	0.852	0.889	0.901	3.71
43) T	Methyl Methacr...			0.545	0.520	0.495	0.430	0.441	0.486	10.21
44) T	2,2,4-Trimethy...			1.935	1.916	1.962	1.708	1.794	1.863	5.79
45) T	t-1,3-Dichloro...			0.514	0.510	0.509	0.459	0.491	0.497	4.60
46) T	cis-1,3-Dichlo...			0.613	0.617	0.637	0.570	0.603	0.608	4.05
47) T	1,1,2-Trichlor...			0.423	0.429	0.420	0.376	0.399	0.410	5.32
48) T	Dibromochlorom...			0.743	0.721	0.746	0.679	0.729	0.723	3.70
49) T	Bromoform			0.622	0.623	0.675	0.616	0.627	0.633	3.82
50) T	4-Methyl-2-Pen...			1.541	1.382	1.435	1.082	1.132	1.314	15.12
51) T	2-Hexanone			1.341	1.311	1.411	0.893	0.938	1.179	20.69
52) T	Tetrachloroethene	0.543	0.452	0.420	0.393	0.391	0.357	0.363	0.417	15.46
53) T	Toluene			1.371	1.295	1.370	1.193	1.244	1.294	6.03
54) T	1,2-Dibromoethane		0.621	0.582	0.605	0.606	0.539	0.562	0.586	5.27
-----ISTD-----										
55) I	Chlorobenzene-d5									
56) T	1,1,1,2-Tetrac...			0.621	0.640	0.635	0.550	0.581	0.605	6.37
57) T	Chlorobenzene			1.129	1.067	1.068	0.956	0.981	1.040	6.79
58) T	Ethyl Benzene			2.029	2.088	2.068	1.832	1.890	1.981	5.73
59) T	m/p-Xylene			1.685	1.682	1.658	1.456	1.500	1.596	6.86
60) T	o-Xylene			1.740	1.697	1.668	1.442	1.490	1.607	8.25
61) T	Styrene			0.792	0.776	0.808	0.726	0.734	0.767	4.68
62) T	Isopropylbenzene			2.517	2.470	2.439	2.170	2.202	2.360	6.83
63) T	1,1,2,2-Tetrac...	1.407	1.176	0.990	0.985	0.968	0.873	0.892	1.042	18.12
64) T	n-propylbenzene			0.645	0.605	0.610	0.542	0.564	0.593	6.81
65) T	tert-Butylbenzene			2.319	2.358	2.315	1.907	1.945	2.169	10.27
66) T	Benzyl Chloride			0.241	0.239	0.258	0.246	0.208	0.238	7.84
67) T	sec-Butylbenzene			3.242	3.219	3.253	2.662	2.709	3.017	10.05
68) S	1-Bromo-4-Fluo...	0.819	0.814	0.824	0.818	0.822	0.831	0.825	0.822	0.69
69) T	p-Isopropyltol...			2.834	2.818	2.811	2.304	2.338	2.621	10.46
70) T	n-Butylbenzene			2.985	2.926	2.953	2.370	2.414	2.730	11.33
71) T	2-Chlorotoluene			1.972	1.924	1.920	1.695	1.744	1.851	6.65
72) T	4-Ethyltoluene			2.055	2.118	2.054	1.765	1.833	1.965	7.92
73) T	1,3,5-Trimethy...			1.894	1.851	1.841	1.517	1.570	1.735	10.19
74) T	1,2,4-Trimethy...			2.348	2.245	2.168	1.671	1.698	2.026	15.71
75) T	1,3-Dichlorobe...			1.119	1.121	1.119	0.948	0.960	1.053	8.64
76) T	1,4-Dichlorobe...			1.138	1.115	1.122	0.944	0.974	1.059	8.69
77) T	1,2-Dichlorobe...			1.085	1.093	1.114	0.904	0.924	1.024	9.88
78) T	Hexachloro-1,3...			1.163	1.136	1.135	0.863	0.874	1.034	14.66
79) T	Naphthalene		2.433	2.395	2.756	2.909	2.004	2.040	2.423	15.11
80) T	Naphthalene,2-...			1.311	1.534	1.777	1.115	1.168	1.381	19.88
81) T	1,2,4-Trichlor...			1.026	1.117	1.224	0.905	0.929	1.040	12.75

(#) = Out of Range

Data Path : Z:\voasrv\HPCHEM1\MSVOA_L\Data\VL041725\
 Data File : VL042349.D
 Acq On : 17 Apr 2025 09:44
 Operator : SY/MD
 Sample : VSTDICCC010
 Misc : 400mL/MSVOA_L
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_L
 ClientSampleId :
 VSTDICCC010

Manual Integrations
 APPROVED

Reviewed By :Semsettin Yesilyurt 04/18/2025
 Supervised By :Mahesh Dadoda 04/18/2025

Quant Time: Apr 18 01:25:08 2025

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_L\methods\VL041725AIR.M

Quant Title : AIR ANALYSIS BY METHOD TO-15 Instrument: MSVOA_L Fri Aug 2025

QLast Update : Fri Apr 18 01:24:29 2025

Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) Bromochloromethane	2.787	49	125455	10.000	ppbv	0.00
33) 1,4-Difluorobenzene	3.959	114	366544	10.000	ppbv	0.00
55) Chlorobenzene-d5	8.882	117	338126	10.000	ppbv	0.00

System Monitoring Compounds

68) 1-Bromo-4-Fluorobenzene 10.377 95 281013 10.114 ppbv 0.00
 Spiked Amount 10.000 Range 65 - 135 Recovery = 101.100%

Target Compounds

						Qvalue
2) Dichlorodifluoromethane	1.499	85	207962	9.382	ppbv	100
3) Chlorodifluoromethane	1.476	51	197334	9.350	ppbv	100
4) Chloromethane	1.531	50	81843	9.330	ppbv	100
5) Vinyl Chloride	1.580	62	80068	8.385	ppbv	100
6) Bromomethane	1.667	94	38128	9.219	ppbv	100
7) Chloroethane	1.703	64	31705	9.182	ppbv	100
8) Dichlorotetrafluoroethane	1.554	85	179696	9.442	ppbv	100
9) Propene	1.483	41	86530	9.002	ppbv	100
10) Heptane	4.978	43	236341	9.138	ppbv	100
11) Trichlorofluoromethane	1.877	101	196596	9.327	ppbv	100
12) 1,1,2-Trichlorotrifluo...	2.136	101	160682	9.323	ppbv	100
13) Ethanol	1.719	45	4439	5.651	ppbv	100
14) Bromoethene	1.780	108	61493	9.466	ppbv	100
15) Acetone	1.835	43	158490	7.717	ppbv	100
16) 1,3-Butadiene	1.609	39	81688	8.470	ppbv	100
17) tert-Butyl alcohol	2.039	59	226280	8.891	ppbv	100
18) 1,1-Dichloroethene	2.033	96	73186	8.833	ppbv	100
19) Isopropyl Alcohol	1.884	45	103029	9.031	ppbv	100
20) Methylene Chloride	2.062	84	62414	7.760	ppbv	100
21) Allyl Chloride	2.094	41	122115	8.611	ppbv	100
22) trans-1,2-Dichloroethene	2.340	96	83278	9.084	ppbv	100
23) Vinyl Acetate	2.460	43	69649	8.378	ppbv	100
24) 1,1-Dichloroethane	2.408	63	161102	9.251	ppbv	100
25) Ethyl Acetate	2.836	43	461315	9.110	ppbv	100
26) Hexane	2.826	57	178691	8.720	ppbv	100
27) Carbon Disulfide	2.149	76	209297	9.539	ppbv	100
28) Methyl tert-Butyl Ether	2.441	73	117924	8.783	ppbv	100
29) Chloroform	2.848	83	269088	9.290	ppbv	100
30) Cyclohexane	3.855	84	157035	9.517	ppbv	100
31) cis-1,2-Dichloroethene	2.719	61	183082	9.397	ppbv	100
32) 1,1,1-Trichloroethane	3.366	97	285336	9.187	ppbv	100
34) 2-Butanone	2.557	43	251704	8.683	ppbv	100
35) Carbon Tetrachloride	3.761	117	282402	8.513	ppbv	100
36) Benzene	3.654	78	367355	9.292	ppbv	100
37) 1,2-Dichloroethane	3.217	62	216999	9.145	ppbv	100
38) Trichloroethene	4.548	130	144767	8.352	ppbv	100
39) 1,2-Dichloropropane	4.311	63	135605	9.197	ppbv	100
40) 1,4-Dioxane	4.580	88	66662	8.348	ppbv #	100
41) Tetrahydrofuran	3.052	42	149905	8.950	ppbv	100
42) Bromodichloromethane	4.489	83	312268	9.456	ppbv	100
43) Methyl Methacrylate	4.881	69	157501	8.876	ppbv	100
44) 2,2,4-Trimethylpentane	4.642	57	626085	9.157	ppbv	100
45) t-1,3-Dichloropropene	6.415	75	168223	9.233	ppbv	100

Data Path : Z:\voasrv\HPCHEM1\MSVOA_L\Data\VL041725\
 Data File : VL042349.D
 Acq On : 17 Apr 2025 09:44
 Operator : SY/MD
 Sample : VSTDICCC010
 Misc : 400mL/MSVOA_L
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_L
 ClientSampleId :
 VSTDICCC010

Manual Integrations
 APPROVED

Reviewed By :Semsettin Yesilyurt 04/18/2025
 Supervised By :Mahesh Dadoda 04/18/2025

Quant Time: Apr 18 01:25:08 2025

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_L\methods\VL041725AIR.M

Quant Title : AIR ANALYSIS BY METHOD TO-15 Instrument: MSVOA_L Fri Aug 2025

QLast Update : Fri Apr 18 01:24:29 2025

Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
46) cis-1,3-Dichloropropene	5.600	75	208811	9.377	ppbv	100
47) 1,1,2-Trichloroethane	6.580	97	137898	9.181	ppbv	100
48) Dibromochloromethane	7.386	129	248954	9.388	ppbv	100
49) Bromoform	9.480	173	225706	9.734	ppbv	100
50) 4-Methyl-2-Pentanone	5.752	43	396548	8.215	ppbv	100
51) 2-Hexanone	7.435	43	327165	7.511	ppbv	100
52) Tetrachloroethene	8.225	164	130720	8.553	ppbv	100
53) Toluene	6.933	91	437258	9.216	ppbv	100
54) 1,2-Dibromoethane	7.652	107	197665	9.235	ppbv	100
56) 1,1,1,2-Tetrachloroethane	8.927	131	186041	9.091	ppbv	100
57) Chlorobenzene	8.920	112	323084	9.187	ppbv	100
58) Ethyl Benzene	9.354	91	619581	9.248	ppbv	100
59) m/p-Xylene	9.552	91	984458m	18.187	ppbv	
60) o-Xylene	9.963	91	487616	8.972	ppbv	100
61) Styrene	9.869	104	245487	9.464	ppbv	100
62) Isopropylbenzene	10.539	105	733788	9.197	ppbv	100
63) 1,1,2,2-Tetrachloroethane	9.963	83	295041	8.355	ppbv	100
64) n-propylbenzene	10.989	120	183390	9.144	ppbv	100
65) tert-Butylbenzene	11.532	119	644953	8.795	ppbv	100
66) Benzyl Chloride	11.617	91	83269	10.333	ppbv	100
67) sec-Butylbenzene	11.769	105	900074	8.824	ppbv	100
69) p-Isopropyltoluene	11.927	119	778987	8.790	ppbv	100
70) n-Butylbenzene	12.283	91	801364	8.682	ppbv	100
71) 2-Chlorotoluene	10.901	91	573039	9.157	ppbv	100
72) 4-Ethyltoluene	11.125	105	596757	8.982	ppbv	100
73) 1,3,5-Trimethylbenzene	11.205	105	512855	8.743	ppbv	100
74) 1,2,4-Trimethylbenzene	11.539	105	565064	8.248	ppbv	100
75) 1,3-Dichlorobenzene	11.607	146	320380	8.997	ppbv	100
76) 1,4-Dichlorobenzene	11.672	146	319089	8.915	ppbv	100
77) 1,2-Dichlorobenzene	11.947	146	305822	8.832	ppbv	100
78) Hexachloro-1,3-Butadiene	13.882	225	291833	8.346	ppbv	100
79) Naphthalene	13.513	128	677578	8.271	ppbv	100
80) Naphthalene,2-methyl-	14.458	142	376995	8.074	ppbv	100
81) 1,2,4-Trichlorobenzene	13.442	180	306119	8.702	ppbv	100

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_L\Data\VL041725\
 Data File : VL042350.D
 Acq On : 17 Apr 2025 10:17
 Operator : SY/MD
 Sample : VSTDICC002
 Misc : 400mL/MSVOA_L
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_L
 ClientSampleId :
 VSTDICC002

Quant Time: Apr 18 01:26:23 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_L\methods\VL041725AIR.M
 Quant Title : AIR ANALYSIS BY METHOD TO-15 Instrument: MSVOA_L Fri Aug 2
 QLast Update : Fri Apr 18 01:24:29 2025
 Response via : Initial Calibration

Manual Integrations APPROVED

Reviewed By : Semsettin Yesilyurt 04/18/2025
 Supervised By : Mahesh Dadoda 04/18/2025

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) Bromochloromethane	2.787	49	123206	10.000	ppbv	0.00
33) 1,4-Difluorobenzene	3.959	114	355918	10.000	ppbv	0.00
55) Chlorobenzene-d5	8.881	117	335603	10.000	ppbv	0.00

System Monitoring Compounds
 68) 1-Bromo-4-Fluorobenzene 10.377 95 275715 9.998 ppbv 0.00
 Spiked Amount 10.000 Range 65 - 135 Recovery = 100.000%

Target Compounds	Qvalue					
2) Dichlorodifluoromethane	1.499	85	45008	2.068	ppbv	95
3) Chlorodifluoromethane	1.476	51	41853	2.019	ppbv	97
4) Chloromethane	1.534	50	17272	2.005	ppbv	93
5) Vinyl Chloride	1.580	62	16979	1.811	ppbv	92
6) Bromomethane	1.667	94	7398	1.821	ppbv #	80
7) Chloroethane	1.706	64	6889	2.031	ppbv	98
8) Dichlorotetrafluoroethane	1.554	85	38955	2.084	ppbv	98
9) Propene	1.483	41	18671	1.978	ppbv	97
10) Heptane	4.978	43	54438m	2.143	ppbv	
11) Trichlorofluoromethane	1.877	101	43035	2.079	ppbv	99
12) 1,1,2-Trichlorotrifluo...	2.140	101	35415	2.092	ppbv	100
13) Ethanol	1.722	45	3034m	3.933	ppbv	
14) Bromoethene	1.780	108	13049	2.045	ppbv #	92
15) Acetone	1.835	43	42949	2.129	ppbv	93
16) 1,3-Butadiene	1.609	39	18110	1.912	ppbv	96
17) tert-Butyl alcohol	2.042	59	51829	2.074	ppbv #	94
18) 1,1-Dichloroethene	2.036	96	15690	1.928	ppbv	87
19) Isopropyl Alcohol	1.887	45	22106	1.973	ppbv #	38
20) Methylene Chloride	2.062	84	14329	1.814	ppbv #	82
21) Allyl Chloride	2.098	41	27340	1.963	ppbv	95
22) trans-1,2-Dichloroethene	2.340	96	18142	2.015	ppbv	97
23) Vinyl Acetate	2.463	43	18454	2.260	ppbv #	91
24) 1,1-Dichloroethane	2.411	63	34077	1.993	ppbv	96
25) Ethyl Acetate	2.835	43	101467	2.040	ppbv	99
26) Hexane	2.826	57	40046	1.990	ppbv	99
27) Carbon Disulfide	2.149	76	43975	2.041	ppbv #	64
28) Methyl tert-Butyl Ether	2.441	73	27905	2.116	ppbv	97
29) Chloroform	2.848	83	58946	2.072	ppbv	100
30) Cyclohexane	3.858	84	33195	2.048	ppbv	94
31) cis-1,2-Dichloroethene	2.722	61	39249	2.051	ppbv	98
32) 1,1,1-Trichloroethane	3.366	97	61442	2.014	ppbv	99
34) 2-Butanone	2.560	43	58249	2.069	ppbv	100
35) Carbon Tetrachloride	3.761	117	60291	1.872	ppbv	88
36) Benzene	3.658	78	79814	2.079	ppbv	98
37) 1,2-Dichloroethane	3.217	62	47398	2.057	ppbv	100
38) Trichloroethene	4.544	130	31826m	1.891	ppbv	
39) 1,2-Dichloropropane	4.315	63	28806	2.012	ppbv	96
40) 1,4-Dioxane	4.593	88	16536m	2.133	ppbv	
41) Tetrahydrofuran	3.059	42	33298	2.047	ppbv	98
42) Bromodichloromethane	4.489	83	64941	2.025	ppbv	97
43) Methyl Methacrylate	4.881	69	35256	2.046	ppbv	97
44) 2,2,4-Trimethylpentane	4.645	57	139641	2.103	ppbv	99
45) t-1,3-Dichloropropene	6.415	75	36265	2.050	ppbv	94

Data Path : Z:\voasrv\HPCHEM1\MSVOA_L\Data\VL041725\
 Data File : VL042350.D
 Acq On : 17 Apr 2025 10:17
 Operator : SY/MD
 Sample : VSTDICC002
 Misc : 400mL/MSVOA_L
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_L
 ClientSampleId :
 VSTDICC002

Manual Integrations
 APPROVED

Reviewed By :Semsettin Yesilyurt 04/18/2025
 Supervised By :Mahesh Dadoda 04/18/2025

Quant Time: Apr 18 01:26:23 2025

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_L\methods\VL041725AIR.M

Quant Title : AIR ANALYSIS BY METHOD TO-15 Instrument: MSVOA_L Fri Aug 2

QLast Update : Fri Apr 18 01:24:29 2025

Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
46) cis-1,3-Dichloropropene	5.600	75	45327m	2.096	ppbv	
47) 1,1,2-Trichloroethane	6.590	97	29879	2.049	ppbv	88
48) Dibromochloromethane	7.389	129	53069	2.061	ppbv	100
49) Bromoform	9.483	173	48069	2.135	ppbv	97
50) 4-Methyl-2-Pentanone	5.758	43	102182	2.180	ppbv	99
51) 2-Hexanone	7.441	43	100443	2.375	ppbv	98
52) Tetrachloroethene	8.228	164	27854	1.877	ppbv	96
53) Toluene	6.933	91	97493	2.116	ppbv	98
54) 1,2-Dibromoethane	7.655	107	43153	2.076	ppbv	87
56) 1,1,1,2-Tetrachloroethane	8.924	131	42601	2.097	ppbv	95
57) Chlorobenzene	8.924	112	71685	2.054	ppbv	98
58) Ethyl Benzene	9.357	91	138810	2.088	ppbv	98
59) m/p-Xylene	9.532	91	222542m	4.142	ppbv	
60) o-Xylene	9.963	91	111972	2.076	ppbv	98
61) Styrene	9.872	104	54229	2.106	ppbv	98
62) Isopropylbenzene	10.539	105	163716	2.067	ppbv	100
63) 1,1,2,2-Tetrachloroethane	9.963	83	64952	1.853	ppbv	99
64) n-propylbenzene	10.989	120	40956	2.058	ppbv	96
65) tert-Butylbenzene	11.532	119	155401	2.135	ppbv	92
66) Benzyl Chloride	11.616	91	17300	2.163	ppbv	98
67) sec-Butylbenzene	11.769	105	218327	2.156	ppbv	100
69) p-Isopropyltoluene	11.927	119	188657	2.145	ppbv	99
70) n-Butylbenzene	12.283	91	198221	2.164	ppbv	99
71) 2-Chlorotoluene	10.901	91	128858	2.075	ppbv	98
72) 4-Ethyltoluene	11.124	105	137872	2.091	ppbv	97
73) 1,3,5-Trimethylbenzene	11.205	105	123580	2.123	ppbv	98
74) 1,2,4-Trimethylbenzene	11.539	105	145510	2.140	ppbv	92
75) 1,3-Dichlorobenzene	11.607	146	75082	2.124	ppbv	98
76) 1,4-Dichlorobenzene	11.671	146	75323	2.120	ppbv	99
77) 1,2-Dichlorobenzene	11.947	146	74786	2.176	ppbv	98
78) Hexachloro-1,3-Butadiene	13.882	225	76158	2.194	ppbv	100
79) Naphthalene	13.513	128	195233	2.401	ppbv	99
80) Naphthalene,2-methyl-	14.458	142	119262	2.574	ppbv	99
81) 1,2,4-Trichlorobenzene	13.442	180	82156	2.353	ppbv	99

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_L\Data\VL041725\
 Data File : VL042351.D
 Acq On : 17 Apr 2025 10:48
 Operator : SY/MD
 Sample : VSTDICC001
 Misc : 400mL/MSVOA_L
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 MSVOA_L
 ClientSampleId :
 VSTDICC001

Quant Time: Apr 18 01:27:33 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_L\methods\VL041725AIR.M
 Quant Title : AIR ANALYSIS BY METHOD TO-15 Instrument: MSVOA_L Fri Aug 2025
 QLast Update : Fri Apr 18 01:24:29 2025
 Response via : Initial Calibration

Manual Integrations APPROVED

Reviewed By : Semsettin Yesilyurt 04/18/2025
 Supervised By : Mahesh Dadoda 04/18/2025

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) Bromochloromethane	2.791	49	124177	10.000	ppbv	0.00
33) 1,4-Difluorobenzene	3.962	114	359907	10.000	ppbv	0.00
55) Chlorobenzene-d5	8.885	117	329090	10.000	ppbv	0.00

System Monitoring Compounds						
68) 1-Bromo-4-Fluorobenzene	10.377	95	269087	9.950	ppbv	0.00
Spiked Amount	10.000	Range	65 - 135	Recovery	=	99.500%

Target Compounds						Qvalue
2) Dichlorodifluoromethane	1.499	85	22816	1.040	ppbv	93
3) Chlorodifluoromethane	1.476	51	20573	0.985	ppbv	97
4) Chloromethane	1.535	50	9008	1.038	ppbv	99
5) Vinyl Chloride	1.580	62	8807	0.932	ppbv	98
6) Bromomethane	1.671	94	4099	1.001	ppbv	91
7) Chloroethane	1.706	64	3559	1.041	ppbv	90
8) Dichlorotetrafluoroethane	1.554	85	19251	1.022	ppbv	100
9) Propene	1.486	41	9892	1.040	ppbv	92
10) Heptane	4.982	43	28507m	1.114	ppbv	
11) Trichlorofluoromethane	1.878	101	22428	1.075	ppbv	99
12) 1,1,2-Trichlorotrifluo...	2.140	101	17865	1.047	ppbv	98
13) Ethanol	1.722	45	1730m	2.225	ppbv	
14) Bromoethene	1.784	108	6766	1.052	ppbv #	89
15) Acetone	1.839	43	23537	1.158	ppbv	97
16) 1,3-Butadiene	1.609	39	10377m	1.087	ppbv	
17) tert-Butyl alcohol	2.043	59	26477	1.051	ppbv #	90
18) 1,1-Dichloroethene	2.036	96	7890	0.962	ppbv	94
19) Isopropyl Alcohol	1.888	45	12175	1.078	ppbv #	39
20) Methylene Chloride	2.062	84	9046	1.136	ppbv	87
21) Allyl Chloride	2.098	41	15057	1.073	ppbv	91
22) trans-1,2-Dichloroethene	2.344	96	9421	1.038	ppbv	96
23) Vinyl Acetate	2.467	43	9234	1.122	ppbv #	90
24) 1,1-Dichloroethane	2.409	63	18236	1.058	ppbv	99
25) Ethyl Acetate	2.842	43	51494	1.027	ppbv #	97
26) Hexane	2.826	57	21350	1.053	ppbv	95
27) Carbon Disulfide	2.153	76	20739	0.955	ppbv #	1
28) Methyl tert-Butyl Ether	2.444	73	14164	1.066	ppbv #	85
29) Chloroform	2.852	83	28973	1.011	ppbv	100
30) Cyclohexane	3.862	84	16547m	1.013	ppbv	
31) cis-1,2-Dichloroethene	2.723	61	19904	1.032	ppbv	97
32) 1,1,1-Trichloroethane	3.376	97	31653	1.030	ppbv	98
34) 2-Butanone	2.564	43	29735	1.045	ppbv	99
35) Carbon Tetrachloride	3.768	117	29875	0.917	ppbv	94
36) Benzene	3.661	78	38707	0.997	ppbv	95
37) 1,2-Dichloroethane	3.224	62	23284	0.999	ppbv	97
38) Trichloroethene	4.551	130	16119	0.947	ppbv	96
39) 1,2-Dichloropropane	4.318	63	14866	1.027	ppbv	96
40) 1,4-Dioxane	4.597	88	8737m	1.114	ppbv	
41) Tetrahydrofuran	3.062	42	17259	1.049	ppbv	96
42) Bromodichloromethane	4.490	83	32721	1.009	ppbv #	98
43) Methyl Methacrylate	4.885	69	18699m	1.073	ppbv	
44) 2,2,4-Trimethylpentane	4.642	57	68946m	1.027	ppbv	
45) t-1,3-Dichloropropene	6.416	75	18347m	1.026	ppbv	

Data Path : Z:\voasrv\HPCHEM1\MSVOA_L\Data\VL041725\
 Data File : VL042351.D
 Acq On : 17 Apr 2025 10:48
 Operator : SY/MD
 Sample : VSTDICC001
 Misc : 400mL/MSVOA_L
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 MSVOA_L
 ClientSampleId :
 VSTDICC001

Manual Integrations
 APPROVED

Reviewed By :Semsettin Yesilyurt 04/18/2025
 Supervised By :Mahesh Dadoda 04/18/2025

Quant Time: Apr 18 01:27:33 2025

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_L\methods\VL041725AIR.M

Quant Title : AIR ANALYSIS BY METHOD TO-15 Instrument: MSVOA_L Fri Aug 2

QLast Update : Fri Apr 18 01:24:29 2025

Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
46) cis-1,3-Dichloropropene	5.603	75	22189m	1.015	ppbv	
47) 1,1,2-Trichloroethane	6.587	97	15457m	1.048	ppbv	
48) Dibromochloromethane	7.390	129	25948	0.997	ppbv	100
49) Bromoform	9.484	173	22413	0.984	ppbv	98
50) 4-Methyl-2-Pentanone	5.765	43	49725	1.049	ppbv	97
51) 2-Hexanone	7.445	43	47178	1.103	ppbv	99
52) Tetrachloroethene	8.228	164	14154	0.943	ppbv	93
53) Toluene	6.940	91	46611	1.001	ppbv	100
54) 1,2-Dibromoethane	7.658	107	21780	1.036	ppbv	93
56) 1,1,1,2-Tetrachloroethane	8.927	131	21052	1.057	ppbv	95
57) Chlorobenzene	8.921	112	35113	1.026	ppbv #	95
58) Ethyl Benzene	9.354	91	68703	1.054	ppbv	96
59) m/p-Xylene	9.552	91	110702m	2.101	ppbv	
60) o-Xylene	9.966	91	55831	1.056	ppbv	99
61) Styrene	9.872	104	25544	1.012	ppbv	100
62) Isopropylbenzene	10.542	105	81286	1.047	ppbv	100
63) 1,1,2,2-Tetrachloroethane	9.966	83	32410	0.943	ppbv	99
64) n-propylbenzene	10.989	120	19898	1.019	ppbv	90
65) tert-Butylbenzene	11.533	119	77598	1.087	ppbv	92
66) Benzyl Chloride	11.617	91	7864	1.003	ppbv #	91
67) sec-Butylbenzene	11.769	105	105919	1.067	ppbv	97
69) p-Isopropyltoluene	11.928	119	92733	1.075	ppbv	97
70) n-Butylbenzene	12.284	91	96306	1.072	ppbv	99
71) 2-Chlorotoluene	10.905	91	63321	1.040	ppbv	99
72) 4-Ethyltoluene	11.128	105	69698	1.078	ppbv	99
73) 1,3,5-Trimethylbenzene	11.206	105	60927	1.067	ppbv	99
74) 1,2,4-Trimethylbenzene	11.539	105	73888	1.108	ppbv	95
75) 1,3-Dichlorobenzene	11.610	146	36877	1.064	ppbv	98
76) 1,4-Dichlorobenzene	11.675	146	36692	1.053	ppbv	98
77) 1,2-Dichlorobenzene	11.950	146	35977	1.068	ppbv	97
78) Hexachloro-1,3-Butadiene	13.886	225	37377	1.098	ppbv	99
79) Naphthalene	13.517	128	90710	1.138	ppbv	99
80) Naphthalene,2-methyl-	14.462	142	50487	1.111	ppbv	98
81) 1,2,4-Trichlorobenzene	13.442	180	36764	1.074	ppbv	99

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

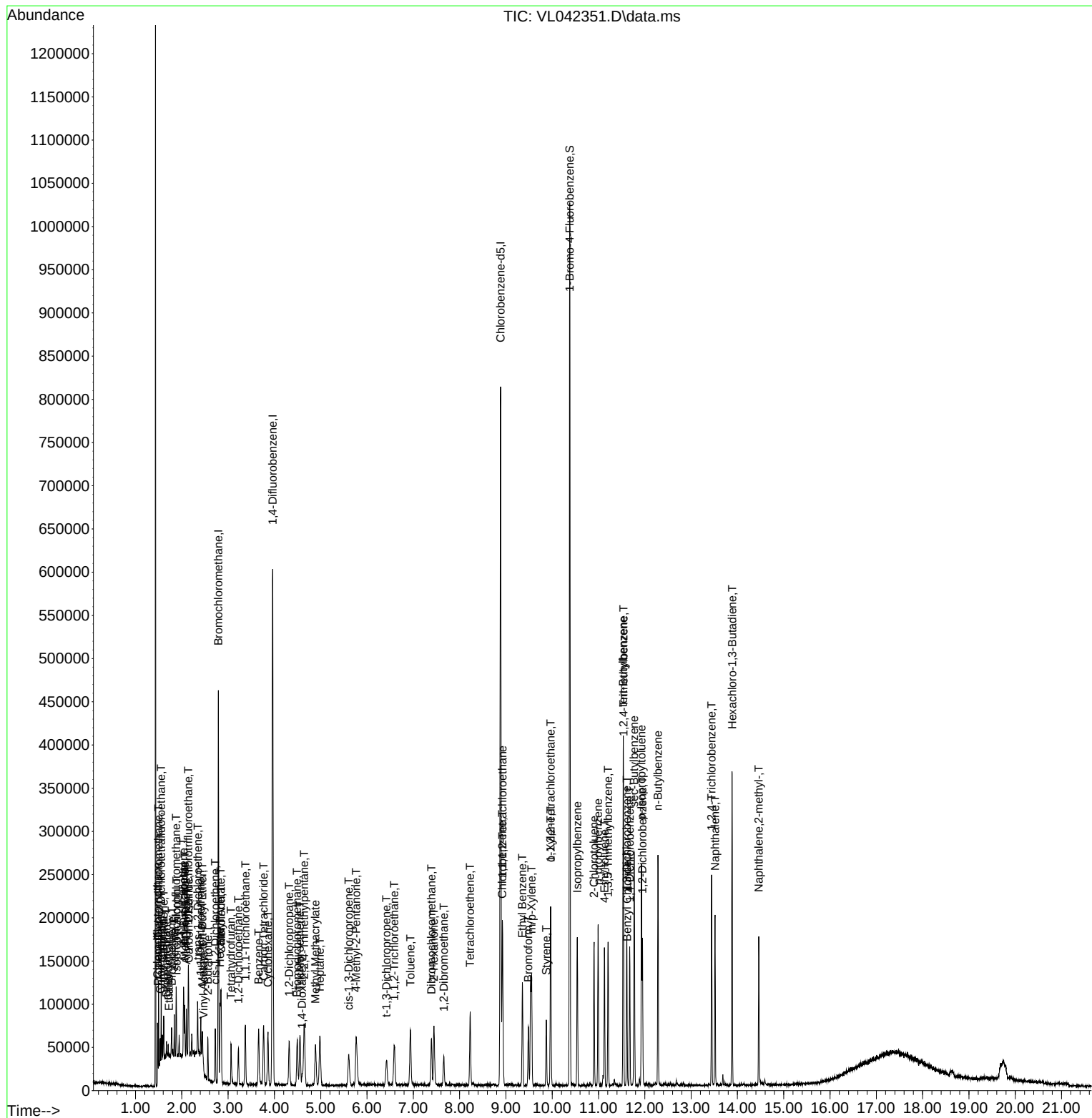
Data Path : Z:\voasrv\HPCHEM1\MSVOA_L\Data\VL041725\
Data File : VL042351.D
Acq On : 17 Apr 2025 10:48
Operator : SY/MD
Sample : VSTDIC001
Misc : 400mL/MSVOA_L
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_L
ClientSampleId :
VSTDIC001

Manual Integrations
APPROVED

Reviewed By :Semsettin Yesilyurt 04/18/2025
Supervised By :Mahesh Dadoda 04/18/2025

Quant Time: Apr 18 01:27:33 2025
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_L\methods\VL041725AIR.M
Quant Title : AIR ANALYSIS BY METHOD TO-15 Instrument: MSVOA_L Fri Aug
QLast Update : Fri Apr 18 01:24:29 2025
Response via : Initial Calibration



Data Path : Z:\voasrv\HPCHEM1\MSVOA_L\Data\VL041725\
 Data File : VL042352.D
 Acq On : 17 Apr 2025 11:19
 Operator : SY/MD
 Sample : VSTDICC0.5
 Misc : 400mL/MSVOA_L
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 MSVOA_L
 ClientSampleId :
 VSTDICC0.5

Quant Time: Apr 18 01:28:45 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_L\methods\VL041725AIR.M
 Quant Title : AIR ANALYSIS BY METHOD TO-15 Instrument: MSVOA_L Fri Aug 2025
 QLast Update : Fri Apr 18 01:24:29 2025
 Response via : Initial Calibration

Manual Integrations APPROVED

Reviewed By : Semsettin Yesilyurt 04/18/2025
 Supervised By : Mahesh Dadoda 04/18/2025

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) Bromochloromethane	2.784	49	120542	10.000	ppbv	0.00
33) 1,4-Difluorobenzene	3.952	114	353137	10.000	ppbv	0.00
55) Chlorobenzene-d5	8.878	117	321644	10.000	ppbv	0.00

System Monitoring Compounds

68) 1-Bromo-4-Fluorobenzene	10.374	95	265161	10.032	ppbv	0.00
Spiked Amount	10.000	Range	65 - 135	Recovery	=	100.300%

Target Compounds					Qvalue	
2) Dichlorodifluoromethane	1.499	85	11036	0.518	ppbv	92
3) Chlorodifluoromethane	1.476	51	11304	0.557	ppbv #	84
4) Chloromethane	1.531	50	4534	0.538	ppbv	97
5) Vinyl Chloride	1.580	62	4406	0.480	ppbv	97
6) Bromomethane	1.667	94	2394	0.602	ppbv	95
7) Chloroethane	1.703	64	1824	0.550	ppbv #	87
8) Dichlorotetrafluoroethane	1.554	85	9381	0.513	ppbv	99
9) Propene	1.483	41	5303	0.574	ppbv	93
10) Heptane	4.978	43	14728m	0.593	ppbv	
11) Trichlorofluoromethane	1.877	101	10150	0.501	ppbv	97
12) 1,1,2-Trichlorotrifluo...	2.136	101	8371	0.506	ppbv	93
13) Ethanol	1.719	45	1656m	2.194	ppbv	
14) Bromoethene	1.780	108	3238	0.519	ppbv #	93
15) Acetone	1.835	43	11859	0.601	ppbv #	74
16) 1,3-Butadiene	1.609	39	6120	0.660	ppbv #	82
17) tert-Butyl alcohol	2.043	59	13640	0.558	ppbv #	85
18) 1,1-Dichloroethene	2.033	96	4430	0.556	ppbv #	87
19) Isopropyl Alcohol	1.884	45	6327	0.577	ppbv #	39
20) Methylene Chloride	2.059	84	5273	0.682	ppbv	93
21) Allyl Chloride	2.094	41	7692	0.565	ppbv #	81
22) trans-1,2-Dichloroethene	2.340	96	4908	0.557	ppbv	92
23) Vinyl Acetate	2.460	43	3953	0.495	ppbv #	57
24) 1,1-Dichloroethane	2.408	63	8753	0.523	ppbv #	94
25) Ethyl Acetate	2.835	43	26686	0.548	ppbv #	93
26) Hexane	2.823	57	11448	0.581	ppbv	91
27) Carbon Disulfide	2.149	76	11472	0.544	ppbv #	1
28) Methyl tert-Butyl Ether	2.441	73	6833	0.530	ppbv #	32
29) Chloroform	2.848	83	14815	0.532	ppbv	93
30) Cyclohexane	3.852	84	8572	0.541	ppbv	94
31) cis-1,2-Dichloroethene	2.722	61	9571	0.511	ppbv	96
32) 1,1,1-Trichloroethane	3.370	97	15413	0.516	ppbv	97
34) 2-Butanone	2.560	43	15640	0.560	ppbv	97
35) Carbon Tetrachloride	3.761	117	14683	0.459	ppbv	87
36) Benzene	3.651	78	20233	0.531	ppbv	97
37) 1,2-Dichloroethane	3.217	62	12194	0.533	ppbv #	95
38) Trichloroethene	4.551	130	7819	0.468	ppbv #	78
39) 1,2-Dichloropropane	4.308	63	7790m	0.548	ppbv	
40) 1,4-Dioxane	4.599	88	4637m	0.603	ppbv	
41) Tetrahydrofuran	3.062	42	9144m	0.567	ppbv	
42) Bromodichloromethane	4.489	83	16642	0.523	ppbv #	91
43) Methyl Methacrylate	4.881	69	9620m	0.563	ppbv	
44) 2,2,4-Trimethylpentane	4.635	57	34170	0.519	ppbv	91
45) t-1,3-Dichloropropene	6.415	75	9073m	0.517	ppbv	

Data Path : Z:\voasrv\HPCHEM1\MSVOA_L\Data\VL041725\
 Data File : VL042352.D
 Acq On : 17 Apr 2025 11:19
 Operator : SY/MD
 Sample : VSTDICC0.5
 Misc : 400mL/MSVOA_L
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 MSVOA_L
 ClientSampleId :
 VSTDICC0.5

Manual Integrations
 APPROVED

Reviewed By :Semsettin Yesilyurt 04/18/2025
 Supervised By :Mahesh Dadoda 04/18/2025

Quant Time: Apr 18 01:28:45 2025

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_L\methods\VL041725AIR.M

Quant Title : AIR ANALYSIS BY METHOD TO-15 Instrument: MSVOA_L Fri Aug 2

QLast Update : Fri Apr 18 01:24:29 2025

Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
46) cis-1,3-Dichloropropene	5.596	75	10827m	0.505	ppbv	
47) 1,1,2-Trichloroethane	6.577	97	7466m	0.516	ppbv	
48) Dibromochloromethane	7.386	129	13122	0.514	ppbv	97
49) Bromoform	9.477	173	10988	0.492	ppbv	99
50) 4-Methyl-2-Pentanone	5.761	43	27206m	0.585	ppbv	
51) 2-Hexanone	7.441	43	23684	0.564	ppbv	98
52) Tetrachloroethene	8.221	164	7412	0.503	ppbv #	88
53) Toluene	6.927	91	24200	0.529	ppbv	97
54) 1,2-Dibromoethane	7.645	107	10282	0.499	ppbv	84
56) 1,1,1,2-Tetrachloroethane	8.930	131	9987	0.513	ppbv #	1
57) Chlorobenzene	8.917	112	18149m	0.543	ppbv	
58) Ethyl Benzene	9.351	91	32624	0.512	ppbv	97
59) m/p-Xylene	9.532	91	54197m	1.053	ppbv	
60) o-Xylene	9.963	91	27983	0.541	ppbv	99
61) Styrene	9.869	104	12736	0.516	ppbv	96
62) Isopropylbenzene	10.539	105	40481	0.533	ppbv	100
63) 1,1,2,2-Tetrachloroethane	9.959	83	15920	0.474	ppbv	99
64) n-propylbenzene	10.985	120	10369	0.544	ppbv	97
65) tert-Butylbenzene	11.529	119	37294	0.535	ppbv	93
66) Benzyl Chloride	11.613	91	3878	0.506	ppbv #	91
67) sec-Butylbenzene	11.765	105	52139	0.537	ppbv	97
69) p-Isopropyltoluene	11.924	119	45574	0.541	ppbv	98
70) n-Butylbenzene	12.280	91	48004	0.547	ppbv	99
71) 2-Chlorotoluene	10.898	91	31709	0.533	ppbv	100
72) 4-Ethyltoluene	11.125	105	33048	0.523	ppbv #	95
73) 1,3,5-Trimethylbenzene	11.202	105	30464	0.546	ppbv	99
74) 1,2,4-Trimethylbenzene	11.536	105	37765	0.579	ppbv #	90
75) 1,3-Dichlorobenzene	11.607	146	18003	0.531	ppbv	99
76) 1,4-Dichlorobenzene	11.668	146	18300	0.537	ppbv	96
77) 1,2-Dichlorobenzene	11.947	146	17446	0.530	ppbv	95
78) Hexachloro-1,3-Butadiene	13.879	225	18705	0.562	ppbv	98
79) Naphthalene	13.513	128	38521	0.494	ppbv #	92
80) Naphthalene,2-methyl-	14.458	142	21076	0.475	ppbv	96
81) 1,2,4-Trichlorobenzene	13.442	180	16504	0.493	ppbv	97

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_L\Data\VL041725\
 Data File : VL042353.D
 Acq On : 17 Apr 2025 11:50
 Operator : SY/MD
 Sample : VSTDICC0.1
 Misc : 400mL/MSVOA_L
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 MSVOA_L
 ClientSampleId :
 VSTDICC0.1

Manual Integrations
 APPROVED

Reviewed By :Semsettin Yesilyurt 04/18/2025
 Supervised By :Mahesh Dadoda 04/18/2025

Quant Time: Apr 18 01:29:39 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_L\methods\VL041725AIR.M
 Quant Title : AIR ANALYSIS BY METHOD T0-15 Instrument: MSVOA_L Fri Aug 2025
 QLast Update : Fri Apr 18 01:24:29 2025
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) Bromochloromethane	2.791	49	119461	10.000	ppbv	0.00
33) 1,4-Difluorobenzene	3.962	114	350484	10.000	ppbv	0.00
55) Chlorobenzene-d5	8.885	117	315004	10.000	ppbv	0.00
System Monitoring Compounds						
68) 1-Bromo-4-Fluorobenzene	10.381	95	256417	9.906	ppbv	0.00
Spiked Amount	10.000	Range	65 - 135	Recovery	=	99.100%
Target Compounds						
					Qvalue	
5) Vinyl Chloride	1.580	62	1088	0.120	ppbv #	74
32) 1,1,1-Trichloroethane	3.367	97	3393m	0.115	ppbv	
35) Carbon Tetrachloride	3.771	117	3204	0.101	ppbv #	94
38) Trichloroethene	4.551	130	1771m	0.107	ppbv	
52) Tetrachloroethene	8.222	164	1583m	0.108	ppbv	
54) 1,2-Dibromoethane	7.662	107	2177m	0.106	ppbv	
63) 1,1,2,2-Tetrachloroethane	9.966	83	3706m	0.113	ppbv	
79) Naphthalene	13.517	128	7664	0.100	ppbv #	92

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

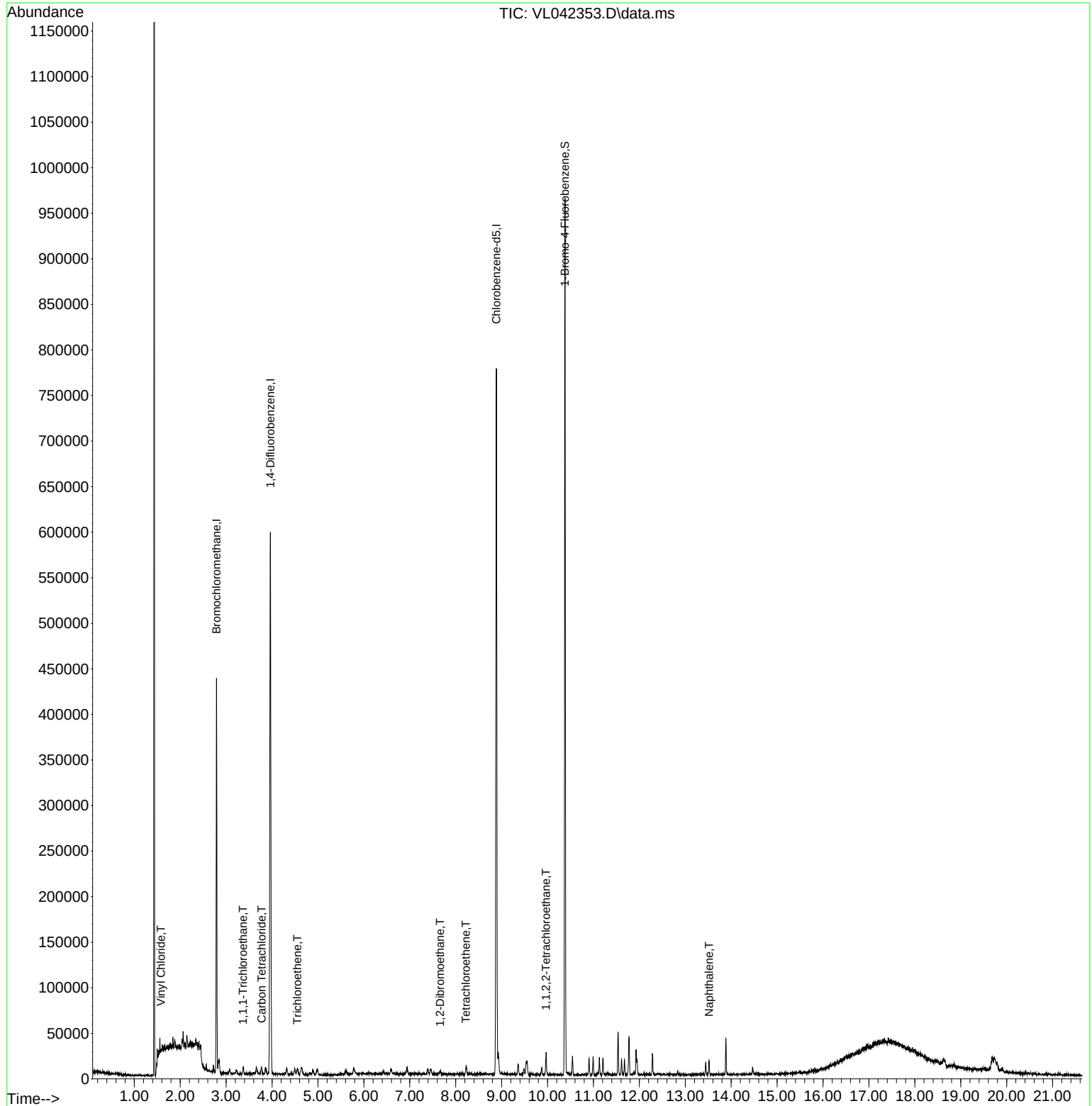
Data Path : Z:\voasrv\HPCHEM1\MSVOA_L\Data\VL041725\
Data File : VL042353.D
Acq On : 17 Apr 2025 11:50
Operator : SY/MD
Sample : VSTDICC0.1
Misc : 400mL/MSVOA_L
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_L
ClientSampleId :
VSTDICC0.1

Manual Integrations
APPROVED

Reviewed By :Semsettin Yesilyurt 04/18/2025
Supervised By :Mahesh Dadoda 04/18/2025

Quant Time: Apr 18 01:29:39 2025
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_L\methods\VL041725AIR.M
Quant Title : AIR ANALYSIS BY METHOD TO-15 Instrument: MSVOA_L Fri Aug
QLast Update : Fri Apr 18 01:24:29 2025
Response via : Initial Calibration



Data Path : Z:\voasrv\HPCHEM1\MSVOA_L\Data\VL041725\
 Data File : VL042354.D
 Acq On : 17 Apr 2025 12:22
 Operator : SY/MD
 Sample : VSTDICC0.03
 Misc : 400mL/MSVOA_L
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 MSVOA_L
 ClientSampleId :
 VSTDICC0.03

Manual Integrations
 APPROVED

Reviewed By :Semsettin Yesilyurt 04/18/2025
 Supervised By :Mahesh Dadoda 04/18/2025

Quant Time: Apr 18 02:07:48 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_L\methods\VL041725AIR.M
 Quant Title : AIR ANALYSIS BY METHOD TO-15 Instrument: MSVOA_L Fri Aug 2
 QLast Update : Fri Apr 18 01:24:29 2025
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) Bromochloromethane	2.790	49	120042	10.000	ppbv	0.00
33) 1,4-Difluorobenzene	3.965	114	349264	10.000	ppbv	0.00
55) Chlorobenzene-d5	8.885	117	312999	10.000	ppbv	0.00
System Monitoring Compounds						
68) 1-Bromo-4-Fluorobenzene	10.380	95	256230	9.962	ppbv	0.00
Spiked Amount	10.000	Range	65 - 135	Recovery	=	99.600%
Target Compounds						
					Qvalue	
5) Vinyl Chloride	1.583	62	361	0.040	ppbv	# 76
32) 1,1,1-Trichloroethane	3.373	97	1120m	0.038	ppbv	
35) Carbon Tetrachloride	3.768	117	1382m	0.044	ppbv	
38) Trichloroethene	4.538	130	674m	0.041	ppbv	
52) Tetrachloroethene	8.228	164	569m	0.039	ppbv	
63) 1,1,2,2-Tetrachloroethane	9.972	83	1321m	0.040	ppbv	

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

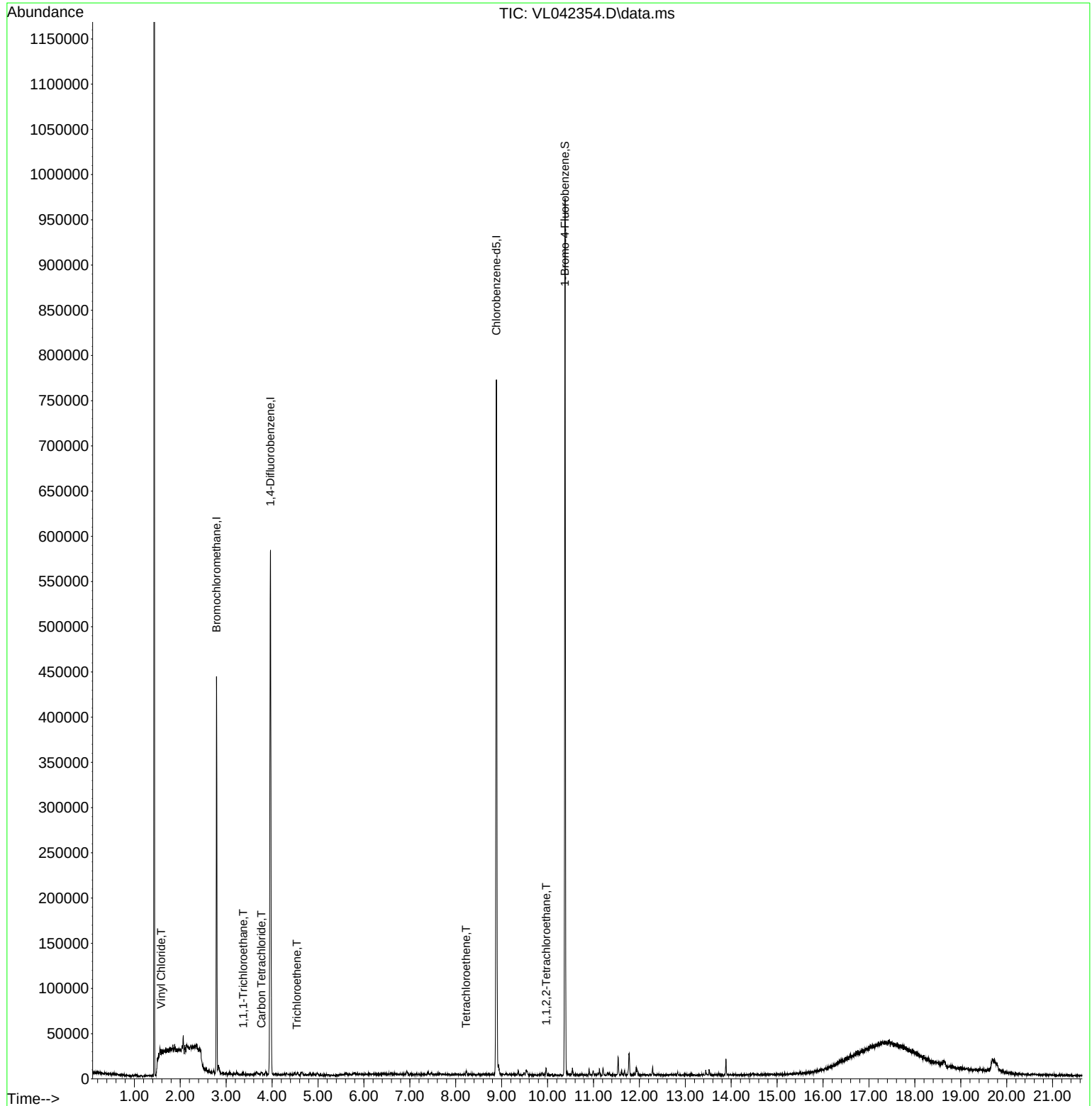
Data Path : Z:\voasrv\HPCHEM1\MSVOA_L\Data\VL041725\
Data File : VL042354.D
Acq On : 17 Apr 2025 12:22
Operator : SY/MD
Sample : VSTDICC0.03
Misc : 400mL/MSVOA_L
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_L
ClientSampleId :
VSTDICC0.03

Manual Integrations
APPROVED

Reviewed By :Semsettin Yesilyurt 04/18/2025
Supervised By :Mahesh Dadoda 04/18/2025

Quant Time: Apr 18 02:07:48 2025
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_L\methods\VL041725AIR.M
Quant Title : AIR ANALYSIS BY METHOD TO-15 Instrument: MSVOA_L Fri Aug
QLast Update : Fri Apr 18 01:24:29 2025
Response via : Initial Calibration



Data Path : Z:\voasrv\HPCHEM1\MSVOA_L\Data\VL041725\
 Data File : VL042355.D
 Acq On : 17 Apr 2025 12:54
 Operator : SY/MD
 Sample : VSTDICC015
 Misc : 400mL/MSVOA_L
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 MSVOA_L
 ClientSampleId :
 VSTDICC015

Manual Integrations
 APPROVED

Reviewed By :Semsettin Yesilyurt 04/18/2025
 Supervised By :Mahesh Dadoda 04/18/2025

Quant Time: Apr 18 01:31:10 2025

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_L\methods\VL041725AIR.M

Quant Title : AIR ANALYSIS BY METHOD TO-15 Instrument: MSVOA_LFri Aug 2025

QLast Update : Fri Apr 18 01:24:29 2025

Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) Bromochloromethane	2.787	49	116071	10.000	ppbv	0.00
33) 1,4-Difluorobenzene	3.959	114	332080	10.000	ppbv	0.00
55) Chlorobenzene-d5	8.882	117	308463	10.000	ppbv	0.00

System Monitoring Compounds

68) 1-Bromo-4-Fluorobenzene 10.374 95 254437 10.038 ppbv 0.00
 Spiked Amount 10.000 Range 65 - 135 Recovery = 100.400%

Target Compounds

						Qvalue
2) Dichlorodifluoromethane	1.499	85	292784	14.277	ppbv	99
3) Chlorodifluoromethane	1.477	51	280042	14.341	ppbv	98
4) Chloromethane	1.532	50	115775	14.266	ppbv	99
5) Vinyl Chloride	1.580	62	112650	12.752	ppbv	100
6) Bromomethane	1.667	94	52805	13.800	ppbv	88
7) Chloroethane	1.703	64	44347	13.881	ppbv	98
8) Dichlorotetrafluoroethane	1.554	85	255029	14.484	ppbv	98
9) Propene	1.483	41	123102	13.842	ppbv	95
10) Heptane	4.979	43	336134	14.047	ppbv	100
11) Trichlorofluoromethane	1.875	101	278059	14.258	ppbv	99
12) 1,1,2-Trichlorotrifluo...	2.137	101	230366	14.447	ppbv	99
13) Ethanol	1.719	45	6324	8.702	ppbv	76
14) Bromoethene	1.781	108	84851	14.117	ppbv	99
15) Acetone	1.833	43	229086	12.056	ppbv	99
16) 1,3-Butadiene	1.609	39	117997	13.224	ppbv	96
17) tert-Butyl alcohol	2.040	59	320605	13.615	ppbv	99
18) 1,1-Dichloroethene	2.033	96	107666	14.045	ppbv	99
19) Isopropyl Alcohol	1.884	45	139004	13.169	ppbv	94
20) Methylene Chloride	2.059	84	91093	12.241	ppbv	90
21) Allyl Chloride	2.095	41	187995	14.329	ppbv	99
22) trans-1,2-Dichloroethene	2.341	96	118492	13.970	ppbv	94
23) Vinyl Acetate	2.460	43	106135	13.799	ppbv #	98
24) 1,1-Dichloroethane	2.409	63	235466	14.615	ppbv	99
25) Ethyl Acetate	2.836	43	663644	14.166	ppbv	100
26) Hexane	2.826	57	260975	13.765	ppbv	97
27) Carbon Disulfide	2.150	76	299098	14.734	ppbv	97
28) Methyl tert-Butyl Ether	2.438	73	174855	14.076	ppbv	95
29) Chloroform	2.849	83	385686	14.393	ppbv	99
30) Cyclohexane	3.855	84	217375	14.238	ppbv	98
31) cis-1,2-Dichloroethene	2.719	61	264972	14.700	ppbv	98
32) 1,1,1-Trichloroethane	3.367	97	412537	14.357	ppbv	99
34) 2-Butanone	2.558	43	367315	13.986	ppbv	100
35) Carbon Tetrachloride	3.762	117	409916	13.639	ppbv	94
36) Benzene	3.655	78	522156	14.578	ppbv	97
37) 1,2-Dichloroethane	3.218	62	319517	14.863	ppbv	100
38) Trichloroethene	4.548	130	208508	13.278	ppbv	98
39) 1,2-Dichloropropane	4.312	63	189697	14.201	ppbv	97
40) 1,4-Dioxane	4.580	88	92114	12.733	ppbv #	95
41) Tetrahydrofuran	3.053	42	214949	14.166	ppbv	98
42) Bromodichloromethane	4.490	83	442735	14.798	ppbv	95
43) Methyl Methacrylate	4.878	69	219757	13.670	ppbv	95
44) 2,2,4-Trimethylpentane	4.642	57	893529	14.425	ppbv	99
45) t-1,3-Dichloropropene	6.412	75	244470	14.810	ppbv	94

Data Path : Z:\voasrv\HPCHEM1\MSVOA_L\Data\VL041725\
 Data File : VL042355.D
 Acq On : 17 Apr 2025 12:54
 Operator : SY/MD
 Sample : VSTDIC015
 Misc : 400mL/MSVOA_L
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 MSVOA_L
 ClientSampleId :
 VSTDIC015

Manual Integrations
 APPROVED

Reviewed By : Semsettin Yesilyurt 04/18/2025
 Supervised By : Mahesh Dadoda 04/18/2025

Quant Time: Apr 18 01:31:10 2025

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_L\methods\VL041725AIR.M

Quant Title : AIR ANALYSIS BY METHOD TO-15 Instrument: MSVOA_L Fri Aug 2

QLast Update : Fri Apr 18 01:24:29 2025

Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
46) cis-1,3-Dichloropropene	5.600	75	300145	14.878	ppbv	90
47) 1,1,2-Trichloroethane	6.581	97	198874	14.615	ppbv	98
48) Dibromochloromethane	7.390	129	362887	15.105	ppbv	98
49) Bromoform	9.484	173	312282	14.865	ppbv	99
50) 4-Methyl-2-Pentanone	5.749	43	563658	12.889	ppbv	99
51) 2-Hexanone	7.435	43	467001	11.834	ppbv #	98
52) Tetrachloroethene	8.225	164	180631	13.046	ppbv	97
53) Toluene	6.933	91	619420	14.411	ppbv	99
54) 1,2-Dibromoethane	7.652	107	280071	14.442	ppbv	92
56) 1,1,1,2-Tetrachloroethane	8.927	131	268626	14.389	ppbv	97
57) Chlorobenzene	8.921	112	454017	14.151	ppbv	97
58) Ethyl Benzene	9.354	91	874424	14.308	ppbv	100
59) m/p-Xylene	9.552	91	1388086m	28.109	ppbv	
60) o-Xylene	9.963	91	689246	13.902	ppbv	99
61) Styrene	9.872	104	339508	14.347	ppbv	99
62) Isopropylbenzene	10.539	105	1018968	13.999	ppbv	100
63) 1,1,2,2-Tetrachloroethane	9.963	83	412898	12.816	ppbv	100
64) n-propylbenzene	10.989	120	260812	14.255	ppbv	94
65) tert-Butylbenzene	11.533	119	899720	13.449	ppbv	93
66) Benzyl Chloride	11.617	91	96016	13.061	ppbv	96
67) sec-Butylbenzene	11.769	105	1253263	13.468	ppbv	100
69) p-Isopropyltoluene	11.928	119	1081850	13.382	ppbv	99
70) n-Butylbenzene	12.284	91	1117042	13.266	ppbv	100
71) 2-Chlorotoluene	10.902	91	806728	14.131	ppbv	100
72) 4-Ethyltoluene	11.125	105	848077	13.992	ppbv	99
73) 1,3,5-Trimethylbenzene	11.206	105	726558	13.578	ppbv	99
74) 1,2,4-Trimethylbenzene	11.539	105	785772	12.572	ppbv	97
75) 1,3-Dichlorobenzene	11.610	146	444086	13.670	ppbv	99
76) 1,4-Dichlorobenzene	11.672	146	450662	13.802	ppbv	98
77) 1,2-Dichlorobenzene	11.947	146	427372	13.529	ppbv	99
78) Hexachloro-1,3-Butadiene	13.883	225	404441	12.679	ppbv	100
79) Naphthalene	13.514	128	944085	12.632	ppbv	98
80) Naphthalene,2-methyl-	14.459	142	540339	12.686	ppbv	99
81) 1,2,4-Trichlorobenzene	13.442	180	429982	13.398	ppbv	98

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

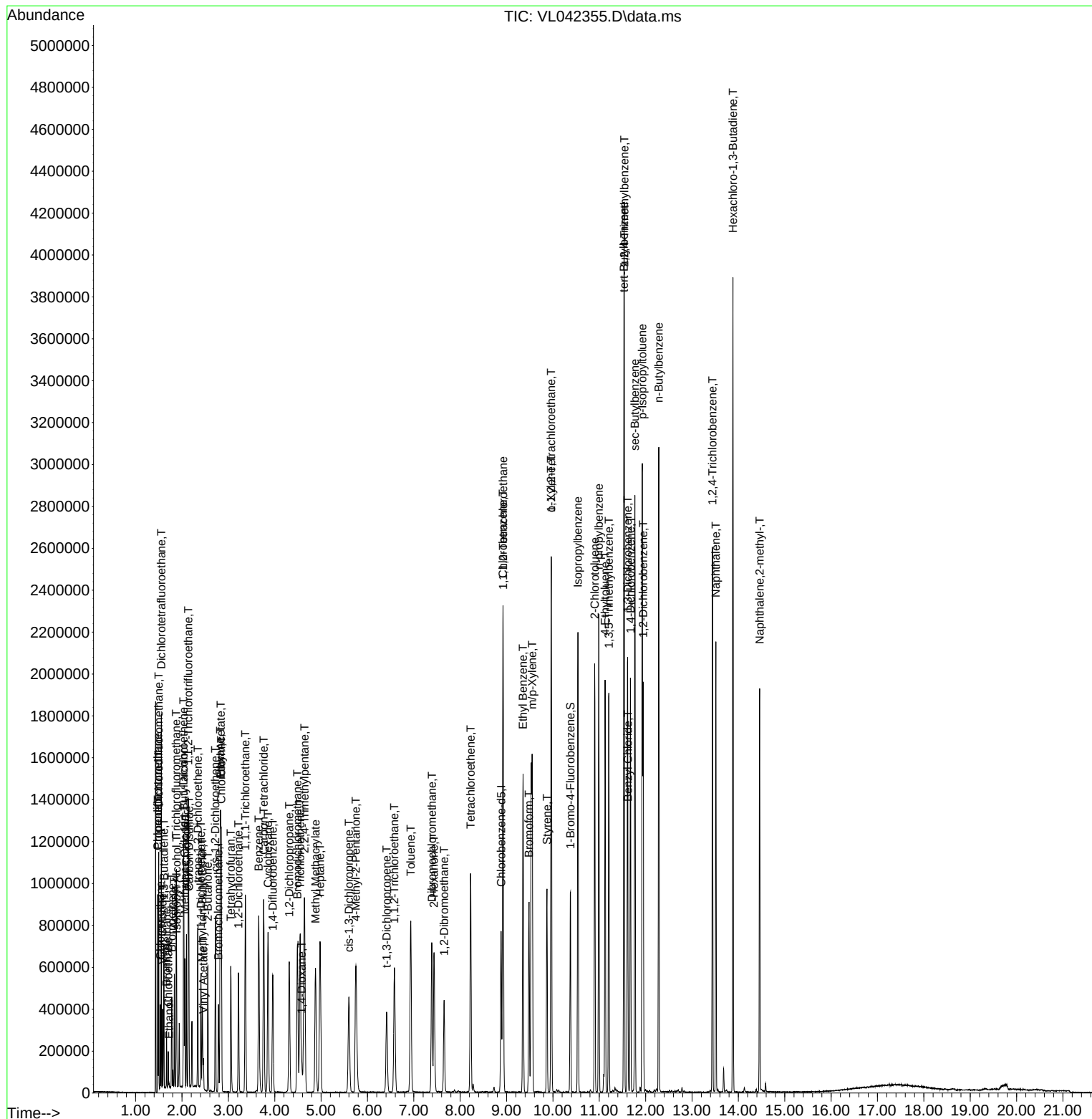
Data Path : Z:\voasrv\HPCHEM1\MSVOA_L\Data\VL041725\
 Data File : VL042355.D
 Acq On : 17 Apr 2025 12:54
 Operator : SY/MD
 Sample : VSTDIC015
 Misc : 400mL/MSVOA_L
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 MSVOA_L
 ClientSampleId :
 VSTDIC015

Manual Integrations APPROVED

Reviewed By :Semsettin Yesilyurt 04/18/2025
 Supervised By :Mahesh Dadoda 04/18/2025

Quant Time: Apr 18 01:31:10 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_L\methods\VL041725AIR.M
 Quant Title : AIR ANALYSIS BY METHOD TO-15 Instrument: MSVOA_L Fri Aug
 QLast Update : Fri Apr 18 01:24:29 2025
 Response via : Initial Calibration



Data Path : Z:\voasrv\HPCHEM1\MSVOA_L\Data\VL041725\
 Data File : VL042356.D
 Acq On : 17 Apr 2025 13:43
 Operator : SY/MD
 Sample : VSTDICV010
 Misc : 400mL/MSVOA_L
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 MSVOA_L
 ClientSampleId :
 ICVVL041725

Manual Integrations
 APPROVED

Reviewed By : Semsettin Yesilyurt 04/18/2025
 Supervised By : Mahesh Dadoda 04/18/2025

Quant Time: Apr 18 02:15:19 2025

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_L\methods\VL041725AIR.M

Quant Title : AIR ANALYSIS BY METHOD TO-15 Instrument: MSVOA_L Fri Aug 2025

QLast Update : Fri Apr 18 02:14:00 2025

Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) Bromochloromethane	2.787	49	113992	10.000	ppbv	0.00
33) 1,4-Difluorobenzene	3.958	114	324528	10.000	ppbv	0.00
55) Chlorobenzene-d5	8.881	117	305357	10.000	ppbv	0.00

System Monitoring Compounds

68) 1-Bromo-4-Fluorobenzene 10.377 95 250149 9.969 ppbv 0.00
 Spiked Amount 10.000 Range 65 - 135 Recovery = 99.700%

Target Compounds

						Qvalue
2) Dichlorodifluoromethane	1.499	85	197623	9.812	ppbv	100
3) Chlorodifluoromethane	1.476	51	191010	9.960	ppbv	99
4) Chloromethane	1.531	50	78080	9.796	ppbv	98
5) Vinyl Chloride	1.580	62	75628	8.717	ppbv	97
6) Bromomethane	1.667	94	35115	9.422	ppbv	84
7) Chloroethane	1.703	64	30158	9.612	ppbv	99
8) Dichlorotetrafluoroethane	1.554	85	172863	9.997	ppbv	98
9) Propene	1.482	41	84636	9.690	ppbv	95
10) Heptane	4.978	43	221840	9.041	ppbv	98
11) Trichlorofluoromethane	1.877	101	187827	9.807	ppbv	99
12) 1,1,2-Trichlorotrifluoro...	2.140	101	148485	9.482	ppbv	100
13) Ethanol	1.722	45	4171	3.005	ppbv	97
14) Bromoethene	1.780	108	58320	9.880	ppbv	99
15) Acetone	1.835	43	158708	8.505	ppbv	100
16) 1,3-Butadiene	1.609	39	78804	8.829	ppbv	95
17) tert-Butyl alcohol	2.039	59	218954	9.468	ppbv	99
18) 1,1-Dichloroethene	2.033	96	67940	9.287	ppbv	94
19) Isopropyl Alcohol	1.887	45	96365	9.296	ppbv	95
20) Methylene Chloride	2.062	84	59203	8.101	ppbv	91
21) Allyl Chloride	2.094	41	120704	9.368	ppbv	98
22) trans-1,2-Dichloroethene	2.340	96	78070	9.373	ppbv	99
23) Vinyl Acetate	2.463	43	73373	9.714	ppbv	99
24) 1,1-Dichloroethane	2.408	63	155571	9.832	ppbv	100
25) Ethyl Acetate	2.835	43	449014	9.759	ppbv	99
26) Hexane	2.826	57	175752	9.439	ppbv	98
27) Carbon Disulfide	2.149	76	191117	9.587	ppbv	100
28) Methyl tert-Butyl Ether	2.441	73	115190	9.442	ppbv	94
29) Chloroform	2.848	83	260855	9.912	ppbv	98
30) Cyclohexane	3.855	84	148246	9.849	ppbv	99
31) cis-1,2-Dichloroethene	2.719	61	179993	10.167	ppbv	99
32) 1,1,1-Trichloroethane	3.369	97	276599	9.336	ppbv	98
34) 2-Butanone	2.557	43	248969	9.700	ppbv	99
35) Carbon Tetrachloride	3.764	117	271862	9.256	ppbv	98
36) Benzene	3.657	78	350545	10.014	ppbv	99
37) 1,2-Dichloroethane	3.217	62	211790	10.081	ppbv	99
38) Trichloroethene	4.551	130	139872	9.143	ppbv	96
39) 1,2-Dichloropropane	4.315	63	129239	9.908	ppbv	99
40) 1,4-Dioxane	4.583	88	61587	8.591	ppbv #	96
41) Tetrahydrofuran	3.052	42	141861	9.479	ppbv	99
42) Bromodichloromethane	4.489	83	296946	10.156	ppbv	96
43) Methyl Methacrylate	4.878	69	149209	9.458	ppbv	96
44) 2,2,4-Trimethylpentane	4.641	57	606376	10.030	ppbv	99
45) t-1,3-Dichloropropene	6.412	75	160750	9.975	ppbv	99

Data Path : Z:\voasrv\HPCHEM1\MSVOA_L\Data\VL041725\
 Data File : VL042356.D
 Acq On : 17 Apr 2025 13:43
 Operator : SY/MD
 Sample : VSTDICV010
 Misc : 400mL/MSVOA_L
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 MSVOA_L
 ClientSampleId :
 ICVVL041725

Manual Integrations
 APPROVED

Reviewed By :Semsettin Yesilyurt 04/18/2025
 Supervised By :Mahesh Dadoda 04/18/2025

Quant Time: Apr 18 02:15:19 2025

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_L\methods\VL041725AIR.M

Quant Title : AIR ANALYSIS BY METHOD TO-15 Instrument: MSVOA_L Fri Aug 2

QLast Update : Fri Apr 18 02:14:00 2025

Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
46) cis-1,3-Dichloropropene	5.599	75	200144	10.148	ppbv	88
47) 1,1,2-Trichloroethane	6.583	97	135575	10.202	ppbv	95
48) Dibromochloromethane	7.389	129	237206	10.103	ppbv	100
49) Bromoform	9.483	173	205630	10.016	ppbv	99
50) 4-Methyl-2-Pentanone	5.748	43	376145	8.819	ppbv	99
51) 2-Hexanone	7.435	43	310521	8.118	ppbv #	99
52) Tetrachloroethene	8.224	164	119424	8.827	ppbv	94
53) Toluene	6.933	91	419462	9.986	ppbv	99
54) 1,2-Dibromoethane	7.652	107	186231	9.792	ppbv	98
56) 1,1,1,2-Tetrachloroethane	8.927	131	182712	9.886	ppbv	97
57) Chlorobenzene	8.920	112	304533	9.589	ppbv	96
58) Ethyl Benzene	9.354	91	583515	9.645	ppbv	99
59) m/p-Xylene	9.532	91	932801m	19.139	ppbv	
60) o-Xylene	9.966	91	464522	9.465	ppbv	99
61) Styrene	9.872	104	229869	9.813	ppbv	98
62) Isopropylbenzene	10.539	105	685965	9.520	ppbv	100
63) 1,1,2,2-Tetrachloroethane	9.962	83	279007	8.773	ppbv	100
64) n-propylbenzene	10.992	120	172654	9.533	ppbv	99
65) tert-Butylbenzene	11.532	119	609500	9.203	ppbv	93
66) Benzyl Chloride	11.613	91	63051	8.664	ppbv	95
67) sec-Butylbenzene	11.769	105	853807	9.268	ppbv	99
69) p-Isopropyltoluene	11.927	119	737861	9.220	ppbv	99
70) n-Butylbenzene	12.283	91	771461	9.255	ppbv	100
71) 2-Chlorotoluene	10.901	91	543794	9.622	ppbv	98
72) 4-Ethyltoluene	11.124	105	563136	9.385	ppbv	100
73) 1,3,5-Trimethylbenzene	11.205	105	486583	9.186	ppbv	99
74) 1,2,4-Trimethylbenzene	11.539	105	531576	8.592	ppbv	97
75) 1,3-Dichlorobenzene	11.610	146	297157	9.240	ppbv	99
76) 1,4-Dichlorobenzene	11.671	146	297697	9.210	ppbv	98
77) 1,2-Dichlorobenzene	11.950	146	289284	9.251	ppbv	100
78) Hexachloro-1,3-Butadiene	13.882	225	276960	8.771	ppbv	100
79) Naphthalene	13.513	128	647349	8.750	ppbv	99
80) Naphthalene,2-methyl-	14.458	142	350259	8.307	ppbv	99
81) 1,2,4-Trichlorobenzene	13.442	180	290260	9.136	ppbv	99

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_L\Data\VL041725\
 Data File : VL042356.D
 Acq On : 17 Apr 2025 13:43
 Operator : SY/MD
 Sample : VSTDICV010
 Misc : 400mL/MSVOA_L
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 MSVOA_L
 ClientSampleId :
 ICSVVL041725

Quant Time: Apr 18 02:15:19 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_L\methods\VL041725AIR.M
 Quant Title : AIR ANALYSIS BY METHOD TO-15 Instrument: MSVOA_L Fri Aug 26 06:05:16 2022
 QLast Update : Fri Apr 18 02:14:00 2025
 Response via : Initial Calibration

Min. RRF : 0.050 Min. Rel. Area : 20% Max. R.T. Dev 0.50min
 Max. RRF Dev : 30% Max. Rel. Area : 150%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
1 I	Bromochloromethane	1.000	1.000	0.0	91	0.00
2 T	Dichlorodifluoromethane	1.767	1.734	1.9	95	0.00
3	Chlorodifluoromethane	1.682	1.676	0.4	97	0.00
4	Chloromethane	0.699	0.685	2.0	95	0.00
5 T	Vinyl Chloride	0.761	0.663	12.9	94	0.00
6 T	Bromomethane	0.327	0.308	5.8	92	0.00
7	Chloroethane	0.275	0.265	3.6	95	0.00
8 T	Dichlorotetrafluoroethane	1.517	1.516	0.1	96	0.00
9 T	Propene	0.766	0.742	3.1	98	0.00
10 T	Heptane	2.153	1.946	9.6	94	0.00
11 T	Trichlorofluoromethane	1.680	1.648	1.9	96	0.00
12 T	1,1,2-Trichlorotrifluoroeth	1.374	1.303	5.2	92	0.00
13	Ethanol	0.122	0.037#	69.7#	94	0.00
14 T	Bromoethene	0.518	0.512	1.2	95	0.00
15 T	Acetone	1.637	1.392	15.0	100	0.00
16 T	1,3-Butadiene	0.783	0.691	11.7	96	0.00
17	tert-Butyl alcohol	2.029	1.921	5.3	97	0.00
18 T	1,1-Dichloroethene	0.642	0.596	7.2	93	0.00
19 T	Isopropyl Alcohol	0.909	0.845	7.0	94	0.00
20 T	Methylene Chloride	0.641	0.519	19.0	95	0.00
21 T	Allyl Chloride	1.130	1.059	6.3	99	0.00
22 T	trans-1,2-Dichloroethene	0.731	0.685	6.3	94	0.00
23 T	Vinyl Acetate	0.663	0.644	2.9	105	0.00
24 T	1,1-Dichloroethane	1.388	1.365	1.7	97	0.00
25 T	Ethyl Acetate	4.036	3.939	2.4	97	0.00
26 T	Hexane	1.633	1.542	5.6	98	0.00
27 T	Carbon Disulfide	1.749	1.677	4.1	91	0.00
28 T	Methyl tert-Butyl Ether	1.070	1.011	5.5	98	0.00
29 T	Chloroform	2.309	2.288	0.9	97	0.00
30 T	Cyclohexane	1.320	1.300	1.5	94	0.00
31 T	cis-1,2-Dichloroethene	1.553	1.579	-1.7	98	0.00
32 T	1,1,1-Trichloroethane	2.599	2.426	6.7	97	0.00
33 I	1,4-Difluorobenzene	1.000	1.000	0.0	89	0.00
34 T	2-Butanone	0.791	0.767	3.0	99	0.00
35 T	Carbon Tetrachloride	0.905	0.838	7.4	96	0.00
36 T	Benzene	1.079	1.080	-0.1	95	0.00
37 T	1,2-Dichloroethane	0.647	0.653	-0.9	98	0.00
38 T	Trichloroethene	0.471	0.431	8.5	97	0.00
39 T	1,2-Dichloropropane	0.402	0.398	1.0	95	0.00
40 T	1,4-Dioxane	0.221	0.190	14.0	92	0.00
41 T	Tetrahydrofuran	0.461	0.437	5.2	95	0.00
42 T	Bromodichloromethane	0.901	0.915	-1.6	95	0.00
43	Methyl Methacrylate	0.486	0.460	5.3	95	0.00
44 T	2,2,4-Trimethylpentane	1.863	1.868	-0.3	97	0.00
45 T	t-1,3-Dichloropropene	0.497	0.495	0.4	96	0.00
46 T	cis-1,3-Dichloropropene	0.608	0.617	-1.5	96	0.00
47 T	1,1,2-Trichloroethane	0.410	0.418	-2.0	98	0.00
48 T	Dibromochloromethane	0.723	0.731	-1.1	95	0.00

Data Path : Z:\voasrv\HPCHEM1\MSVOA_L\Data\VL041725\
 Data File : VL042356.D
 Acq On : 17 Apr 2025 13:43
 Operator : SY/MD
 Sample : VSTDICV010
 Misc : 400mL/MSVOA_L
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 MSVOA_L
 ClientSampleId :
 ICVVL041725

Quant Time: Apr 18 02:15:19 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_L\methods\VL041725AIR.M
 Quant Title : AIR ANALYSIS BY METHOD TO-15 Instrument: MSVOA_L Fri Aug 26 06:05:16 2022
 QLast Update : Fri Apr 18 02:14:00 2025
 Response via : Initial Calibration

Min. RRF : 0.050 Min. Rel. Area : 20% Max. R.T. Dev 0.50min
 Max. RRF Dev : 30% Max. Rel. Area : 150%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
49 T	Bromoform	0.633	0.634	-0.2	91	0.00
50 T	4-Methyl-2-Pentanone	1.314	1.159	11.8	95	0.00
51 T	2-Hexanone	1.179	0.957	18.8	95	0.00
52 T	Tetrachloroethene	0.417	0.368	11.8	91	0.00
53 T	Toluene	1.294	1.293	0.1	96	0.00
54 T	1,2-Dibromoethane	0.586	0.574	2.0	94	0.00
55 I	Chlorobenzene-d5	1.000	1.000	0.0	90	0.00
56	1,1,1,2-Tetrachloroethane	0.605	0.598	1.2	98	0.00
57 T	Chlorobenzene	1.040	0.997	4.1	94	0.00
58 T	Ethyl Benzene	1.981	1.911	3.5	94	0.00
59 T	m/p-Xylene	1.596	1.527	4.3	95	-0.02
60 T	o-Xylene	1.607	1.521	5.4	95	0.00
61 T	Styrene	0.767	0.753	1.8	94	0.00
62	Isopropylbenzene	2.360	2.246	4.8	93	0.00
63 T	1,1,2,2-Tetrachloroethane	1.042	0.914	12.3	95	0.00
64	n-propylbenzene	0.593	0.565	4.7	94	0.00
65	tert-Butylbenzene	2.169	1.996	8.0	95	0.00
66 T	Benzyl Chloride	0.238	0.206	13.4	76	0.00
67	sec-Butylbenzene	3.017	2.796	7.3	95	0.00
68 S	1-Bromo-4-Fluorobenzene	0.822	0.819	0.4	89	0.00
69	p-Isopropyltoluene	2.621	2.416	7.8	95	0.00
70	n-Butylbenzene	2.730	2.526	7.5	96	0.00
71	2-Chlorotoluene	1.851	1.781	3.8	95	0.00
72 T	4-Ethyltoluene	1.965	1.844	6.2	94	0.00
73 T	1,3,5-Trimethylbenzene	1.735	1.593	8.2	95	0.00
74 T	1,2,4-Trimethylbenzene	2.026	1.741	14.1	94	0.00
75 T	1,3-Dichlorobenzene	1.053	0.973	7.6	93	0.00
76 T	1,4-Dichlorobenzene	1.059	0.975	7.9	93	0.00
77 T	1,2-Dichlorobenzene	1.024	0.947	7.5	95	0.00
78 T	Hexachloro-1,3-Butadiene	1.034	0.907	12.3	95	0.00
79 T	Naphthalene	2.423	2.120	12.5	96	0.00
80 T	Naphthalene,2-methyl-	1.381	1.147	16.9	93	0.00
81 T	1,2,4-Trichlorobenzene	1.040	0.951	8.6	95	0.00

(#) = Out of Range

SPCC's out = 0 CCC's out = 0

VOLATILE CONTINUING CALIBRATION CHECK

Lab Name: CHEMTECH Contract: GFEL01

Lab Code: CHEM Case No.: Q1801 SAS No.: Q1801 SDG No.: Q1801

Instrument ID: MSVOA_L Calibration Date/Time: 04/21/2025 08:41

Lab File ID: VL042395.D Init. Calib. Date(s): 04/17/2025 04/17/2025

Heated Purge: (Y/N) N Init. Calib. Time(s): 09:44 12:54

GC Column: RTX-1 ID: 0.32 (mm)

COMPOUND	RRF	RRF010	MIN RRF	%D	MAX%D
Dichlorodifluoromethane	1.767	1.812		2.55	30
Chloromethane	0.699	0.661		-5.44	30
Vinyl Chloride	0.761	0.665		-12.61	30
Bromomethane	0.327	0.310		-5.2	30
Chloroethane	0.275	0.270		-1.82	30
Tetrahydrofuran	0.461	0.355		-22.99	30
Trichlorofluoromethane	1.680	1.735		3.27	30
1,1,2-Trichlorotrifluoroethane	1.374	1.307		-4.88	30
Dichlorotetrafluoroethane	1.517	1.541		1.58	30
tert-Butyl alcohol	2.029	1.870		-7.84	30
Heptane	2.153	1.512		-29.77	30
1,1-Dichloroethene	0.642	0.593		-7.63	30
Acetone	1.637	1.420		-13.26	30
Carbon Disulfide	1.749	1.689		-3.43	30
Methyl tert-Butyl Ether	1.070	0.900		-15.89	30
Methylene Chloride	0.641	0.507		-20.91	30
trans-1,2-Dichloroethene	0.731	0.672		-8.07	30
1,1-Dichloroethane	1.388	1.305		-5.98	30
Cyclohexane	1.320	1.069		-19.01	30
2-Butanone	0.791	0.658		-16.81	30
Carbon Tetrachloride	0.905	0.870		-3.87	30
cis-1,2-Dichloroethene	1.553	1.406		-9.47	30
Chloroform	2.309	2.194		-4.98	30
1,1,1-Trichloroethane	2.599	2.413		-7.16	30
2,2,4-Trimethylpentane	1.863	1.516		-18.63	30
Benzene	1.079	0.908		-15.85	30
1,2-Dichloroethane	0.647	0.661		2.16	30
Trichloroethene	0.471	0.368		-21.87	30
1,2-Dichloropropane	0.402	0.322		-19.9	30
Bromodichloromethane	0.901	0.883		-2	30
4-Methyl-2-Pentanone	1.314	0.961		-26.86	30
Toluene	1.294	1.062		-17.93	30
t-1,3-Dichloropropene	0.497	0.427		-14.09	30
cis-1,3-Dichloropropene	0.608	0.522		-14.15	30
1,1,2-Trichloroethane	0.410	0.357		-12.93	30
Dibromochloromethane	0.723	0.686		-5.12	30
1,2-Dibromoethane	0.586	0.507		-13.48	30
Tetrachloroethene	0.417	0.312		-25.18	30

All other compounds must meet a minimum RRF of 0.010.
RRF of 1,4-Dioxane = Value should be divide by 1000.

VOLATILE CONTINUING CALIBRATION CHECK

Lab Name: CHEMTECH Contract: GFEL01

Lab Code: CHEM Case No.: Q1801 SAS No.: Q1801 SDG No.: Q1801

Instrument ID: MSVOA_L Calibration Date/Time: 04/21/2025 08:41

Lab File ID: VL042395.D Init. Calib. Date(s): 04/17/2025 04/17/2025

Heated Purge: (Y/N) N Init. Calib. Time(s): 09:44 12:54

GC Column: RTX-1 ID: 0.32 (mm)

COMPOUND	RRF	RRF010	MIN RRF	%D	MAX%D
Chlorobenzene	1.040	0.930		-10.58	30
Ethyl Benzene	1.981	1.743		-12.01	30
m/p-Xylene	1.596	1.426		-10.65	30
o-Xylene	1.607	1.459		-9.21	30
Styrene	0.767	0.649		-15.39	30
Bromoform	0.633	0.559		-11.69	30
1,1,2,2-Tetrachloroethane	1.042	0.842		-19.19	30
2-Chlorotoluene	1.851	1.660		-10.32	30
1,3,5-Trimethylbenzene	1.735	1.518		-12.51	30
1,2,4-Trimethylbenzene	2.026	1.690		-16.58	30
1,3-Dichlorobenzene	1.053	0.888		-15.67	30
1,4-Dichlorobenzene	1.059	0.894		-15.58	30
1,2-Dichlorobenzene	1.024	0.875		-14.55	30
1,2,4-Trichlorobenzene	1.040	0.844		-18.85	30
Hexachloro-1,3-Butadiene	1.034	0.828		-19.92	30
1,3-Butadiene	0.783	0.686		-12.39	30
Naphthalene	2.423	1.920		-20.76	30
4-Ethyltoluene	1.965	1.755		-10.69	30
1-Bromo-4-Fluorobenzene	0.822	0.882		7.3	30
Hexane	1.633	1.272		-22.11	30
Allyl Chloride	1.130	1.024		-9.38	30
1,4-Dioxane	220.893	160.950		-27.14	30
Methyl Methacrylate	0.486	0.373		-23.25	30

All other compounds must meet a minimum RRF of 0.010.
RRF of 1,4-Dioxane = Value should be divide by 1000.

Data Path : Z:\voasrv\HPCHEM1\MSVOA_L\Data\VL042125\
 Data File : VL042395.D
 Acq On : 21 Apr 2025 08:41
 Operator : SY/MD
 Sample : VSTDCCC010
 Misc : 400mL/MSVOA_L
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_L
 ClientSampleId :
 VSTDCCC010

Manual Integrations
 APPROVED

Reviewed By : Semsettin Yesilyurt 04/22/2025
 Supervised By : Mahesh Dadoda 04/22/2025

Quant Time: Apr 22 02:50:13 2025

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_L\methods\VL041725AIR.M

Quant Title : AIR ANALYSIS BY METHOD TO-15 Instrument: MSVOA_L Fri Aug 2

QLast Update : Fri Apr 18 02:14:00 2025

Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) Bromochloromethane	2.784	49	118349	10.000	ppbv	0.00
33) 1,4-Difluorobenzene	3.952	114	334726	10.000	ppbv	0.00
55) Chlorobenzene-d5	8.875	117	297161	10.000	ppbv	0.00

System Monitoring Compounds

68) 1-Bromo-4-Fluorobenzene 10.370 95 262155 10.736 ppbv 0.00
 Spiked Amount 10.000 Range 65 - 135 Recovery = 107.400%

Target Compounds

						Qvalue
2) Dichlorodifluoromethane	1.496	85	214405	10.253	ppbv	95
3) Chlorodifluoromethane	1.473	51	172276	8.652	ppbv	96
4) Chloromethane	1.531	50	78242	9.455	ppbv	99
5) Vinyl Chloride	1.576	62	78692	8.736	ppbv	98
6) Bromomethane	1.664	94	36698	9.484	ppbv	92
7) Chloroethane	1.703	64	31945	9.806	ppbv	98
8) Dichlorotetrafluoroethane	1.551	85	182317	10.155	ppbv	94
9) Propene	1.483	41	67297	7.422	ppbv	98
10) Heptane	4.972	43	178981	7.026	ppbv	98
11) Trichlorofluoromethane	1.874	101	205368	10.328	ppbv	98
12) 1,1,2-Trichlorotrifluo...	2.136	101	154701	9.515	ppbv	99
13) Ethanol	1.719	45	4511	3.130	ppbv	88
14) Bromoethene	1.780	108	59325	9.680	ppbv	98
15) Acetone	1.832	43	168061	8.675	ppbv	99
16) 1,3-Butadiene	1.606	39	81189	8.762	ppbv	96
17) tert-Butyl alcohol	2.036	59	221320	9.218	ppbv	100
18) 1,1-Dichloroethene	2.030	96	70214	9.244	ppbv	97
19) Isopropyl Alcohol	1.884	45	98218	9.126	ppbv	94
20) Methylene Chloride	2.059	84	60052	7.915	ppbv	92
21) Allyl Chloride	2.091	41	121139	9.056	ppbv	97
22) trans-1,2-Dichloroethene	2.337	96	79496	9.192	ppbv	92
23) Vinyl Acetate	2.457	43	57298	7.306	ppbv	97
24) 1,1-Dichloroethane	2.405	63	154500	9.405	ppbv	100
25) Ethyl Acetate	2.832	43	392210	8.211	ppbv	99
26) Hexane	2.823	57	150573	7.789	ppbv	98
27) Carbon Disulfide	2.146	76	199908	9.658	ppbv	98
28) Methyl tert-Butyl Ether	2.437	73	106569	8.414	ppbv	95
29) Chloroform	2.845	83	259635	9.502	ppbv	98
30) Cyclohexane	3.852	84	126496	8.095	ppbv	97
31) cis-1,2-Dichloroethene	2.716	61	166357	9.051	ppbv	98
32) 1,1,1-Trichloroethane	3.363	97	285590	9.284	ppbv	99
34) 2-Butanone	2.554	43	220144	8.316	ppbv	99
35) Carbon Tetrachloride	3.755	117	291229	9.614	ppbv	91
36) Benzene	3.651	78	303987	8.420	ppbv	97
37) 1,2-Dichloroethane	3.214	62	221151	10.206	ppbv	97
38) Trichloroethene	4.541	130	123150	7.804	ppbv	94
39) 1,2-Dichloropropane	4.305	63	107877	8.018	ppbv #	88
40) 1,4-Dioxane	4.574	88	53874	7.286	ppbv #	94
41) Tetrahydrofuran	3.049	42	118863	7.701	ppbv	99
42) Bromodichloromethane	4.483	83	295552	9.800	ppbv	98
43) Methyl Methacrylate	4.875	69	124747	7.667	ppbv	95
44) 2,2,4-Trimethylpentane	4.629	57	507437	8.138	ppbv	95
45) t-1,3-Dichloropropene	6.406	75	142879	8.596	ppbv	96

Data Path : Z:\voasrv\HPCHEM1\MSVOA_L\Data\VL042125\
 Data File : VL042395.D
 Acq On : 21 Apr 2025 08:41
 Operator : SY/MD
 Sample : VSTDCCC010
 Misc : 400mL/MSVOA_L
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_L
 ClientSampleId :
 VSTDCCC010

Manual Integrations
 APPROVED

Reviewed By :Semsettin Yesilyurt 04/22/2025
 Supervised By :Mahesh Dadoda 04/22/2025

Quant Time: Apr 22 02:50:13 2025

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_L\methods\VL041725AIR.M

Quant Title : AIR ANALYSIS BY METHOD TO-15 Instrument: MSVOA_L Fri Aug 2

QLast Update : Fri Apr 18 02:14:00 2025

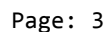
Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
46) cis-1,3-Dichloropropene	5.590	75	174844	8.595	ppbv	88
47) 1,1,2-Trichloroethane	6.571	97	119399	8.711	ppbv	92
48) Dibromochloromethane	7.380	129	229466	9.476	ppbv	99
49) Bromoform	9.480	173	187094	8.836	ppbv	99
50) 4-Methyl-2-Pentanone	5.742	43	321752	7.314	ppbv	99
51) 2-Hexanone	7.425	43	260009	6.590	ppbv #	95
52) Tetrachloroethene	8.218	164	104294	7.474	ppbv	90
53) Toluene	6.923	91	355540	8.206	ppbv	99
54) 1,2-Dibromoethane	7.642	107	169774	8.654	ppbv	91
56) 1,1,1,2-Tetrachloroethane	8.917	131	170285	9.468	ppbv	96
57) Chlorobenzene	8.917	112	276428	8.944	ppbv	98
58) Ethyl Benzene	9.348	91	517937	8.797	ppbv	100
59) m/p-Xylene	9.545	91	847455m	17.868	ppbv	
60) o-Xylene	9.959	91	433569	9.078	ppbv	96
61) Styrene	9.866	104	192729	8.454	ppbv	96
62) Isopropylbenzene	10.535	105	625514	8.920	ppbv	98
63) 1,1,2,2-Tetrachloroethane	9.956	83	250238	8.085	ppbv	100
64) n-propylbenzene	10.985	120	154114	8.744	ppbv	97
65) tert-Butylbenzene	11.526	119	570548	8.853	ppbv	96
66) Benzyl Chloride	11.610	91	57090	8.061	ppbv	99
67) sec-Butylbenzene	11.765	105	797738	8.899	ppbv	99
69) p-Isopropyltoluene	11.921	119	684482	8.789	ppbv	97
70) n-Butylbenzene	12.277	91	720062	8.877	ppbv	97
71) 2-Chlorotoluene	10.898	91	493192	8.967	ppbv	97
72) 4-Ethyltoluene	11.121	105	521390	8.929	ppbv	97
73) 1,3,5-Trimethylbenzene	11.199	105	451158	8.752	ppbv	96
74) 1,2,4-Trimethylbenzene	11.532	105	502117	8.340	ppbv #	87
75) 1,3-Dichlorobenzene	11.604	146	264024	8.436	ppbv	97
76) 1,4-Dichlorobenzene	11.668	146	265655	8.445	ppbv	98
77) 1,2-Dichlorobenzene	11.943	146	260133	8.548	ppbv	97
78) Hexachloro-1,3-Butadiene	13.879	225	245910	8.002	ppbv	99
79) Naphthalene	13.510	128	570593	7.925	ppbv	100
80) Naphthalene,2-methyl-	14.455	142	304139	7.412	ppbv	96
81) 1,2,4-Trichlorobenzene	13.439	180	250801	8.112	ppbv	99

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

Manual Integrations APPROVED

Quant Time: Apr 22 02:50:13 2025
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_L\methods\VL041725AIR.M
Quant Title : AIR ANALYSIS BY METHOD TO-15 Instrument: MSVOA_LFri Aug
QLast Update : Fri Apr 18 02:14:00 2025
Response via : Initial Calibration



Data Path : Z:\voasrv\HPCHEM1\MSVOA_L\Data\VL042125\
 Data File : VL042395.D
 Acq On : 21 Apr 2025 08:41
 Operator : SY/MD
 Sample : VSTDCCC010
 Misc : 400mL/MSVOA_L
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_L
 LabSampleId :
 VSTDCCC010

Quant Time: Apr 22 02:50:13 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_L\methods\VL041725AIR.M
 Quant Title : AIR ANALYSIS BY METHOD TO-15 Instrument: MSVOA_L Fri Aug 26 06:05:16 2022
 QLast Update : Fri Apr 18 02:14:00 2025
 Response via : Initial Calibration

Min. RRF : 0.050 Min. Rel. Area : 20% Max. R.T. Dev 0.50min
 Max. RRF Dev : 30% Max. Rel. Area : 150%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
1 I	Bromochloromethane	1.000	1.000	0.0	94	0.00
2 T	Dichlorodifluoromethane	1.767	1.812	-2.5	103	0.00
3	Chlorodifluoromethane	1.682	1.456	13.4	87	0.00
4	Chloromethane	0.699	0.661	5.4	96	0.00
5 T	Vinyl Chloride	0.761	0.665	12.6	98	0.00
6 T	Bromomethane	0.327	0.310	5.2	96	0.00
7	Chloroethane	0.275	0.270	1.8	101	0.00
8 T	Dichlorotetrafluoroethane	1.517	1.541	-1.6	101	0.00
9 T	Propene	0.766	0.569	25.7	78	0.00
10 T	Heptane	2.153	1.512	29.8	76	0.00
11 T	Trichlorofluoromethane	1.680	1.735	-3.3	104	0.00
12 T	1,1,2-Trichlorotrifluoroeth	1.374	1.307	4.9	96	0.00
13	Ethanol	0.122	0.038#	68.9#	102	0.00
14 T	Bromoethene	0.518	0.501	3.3	96	0.00
15 T	Acetone	1.637	1.420	13.3	106	0.00
16 T	1,3-Butadiene	0.783	0.686	12.4	99	0.00
17	tert-Butyl alcohol	2.029	1.870	7.8	98	0.00
18 T	1,1-Dichloroethene	0.642	0.593	7.6	96	0.00
19 T	Isopropyl Alcohol	0.909	0.830	8.7	95	0.00
20 T	Methylene Chloride	0.641	0.507	20.9	96	0.00
21 T	Allyl Chloride	1.130	1.024	9.4	99	0.00
22 T	trans-1,2-Dichloroethene	0.731	0.672	8.1	95	0.00
23 T	Vinyl Acetate	0.663	0.484	27.0	82	0.00
24 T	1,1-Dichloroethane	1.388	1.305	6.0	96	0.00
25 T	Ethyl Acetate	4.036	3.314	17.9	85	0.00
26 T	Hexane	1.633	1.272	22.1	84	0.00
27 T	Carbon Disulfide	1.749	1.689	3.4	96	0.00
28 T	Methyl tert-Butyl Ether	1.070	0.900	15.9	90	0.00
29 T	Chloroform	2.309	2.194	5.0	96	0.00
30 T	Cyclohexane	1.320	1.069	19.0	81	0.00
31 T	cis-1,2-Dichloroethene	1.553	1.406	9.5	91	0.00
32 T	1,1,1-Trichloroethane	2.599	2.413	7.2	100	0.00
33 I	1,4-Difluorobenzene	1.000	1.000	0.0	91	0.00
34 T	2-Butanone	0.791	0.658	16.8	87	0.00
35 T	Carbon Tetrachloride	0.905	0.870	3.9	103	0.00
36 T	Benzene	1.079	0.908	15.8	83	0.00
37 T	1,2-Dichloroethane	0.647	0.661	-2.2	102	0.00
38 T	Trichloroethene	0.471	0.368	21.9	85	0.00
39 T	1,2-Dichloropropane	0.402	0.322	19.9	80	0.00
40 T	1,4-Dioxane	0.221	0.161	27.1	81	0.00
41 T	Tetrahydrofuran	0.461	0.355	23.0	79	0.00
42 T	Bromodichloromethane	0.901	0.883	2.0	95	0.00
43	Methyl Methacrylate	0.486	0.373	23.3	79	0.00
44 T	2,2,4-Trimethylpentane	1.863	1.516	18.6	81	-0.01
45 T	t-1,3-Dichloropropene	0.497	0.427	14.1	85	0.00
46 T	cis-1,3-Dichloropropene	0.608	0.522	14.1	84	0.00
47 T	1,1,2-Trichloroethane	0.410	0.357	12.9	87	0.00
48 T	Dibromochloromethane	0.723	0.686	5.1	92	0.00

Data Path : Z:\voasrv\HPCHEM1\MSVOA_L\Data\VL042125\
 Data File : VL042395.D
 Acq On : 21 Apr 2025 08:41
 Operator : SY/MD
 Sample : VSTDCCC010
 Misc : 400mL/MSVOA_L
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_L
 LabSampleId :
 VSTDCCC010

Quant Time: Apr 22 02:50:13 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_L\methods\VL041725AIR.M
 Quant Title : AIR ANALYSIS BY METHOD TO-15 Instrument: MSVOA_L Fri Aug 26 06:05:16 2022
 QLast Update : Fri Apr 18 02:14:00 2025
 Response via : Initial Calibration

Min. RRF : 0.050 Min. Rel. Area : 20% Max. R.T. Dev 0.50min
 Max. RRF Dev : 30% Max. Rel. Area : 150%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
49 T	Bromoform	0.633	0.559	11.7	83	0.00
50 T	4-Methyl-2-Pentanone	1.314	0.961	26.9	81	0.00
51 T	2-Hexanone	1.179	0.777	34.1#	79	0.00
52 T	Tetrachloroethene	0.417	0.312	25.2	80	0.00
53 T	Toluene	1.294	1.062	17.9	81	0.00
54 T	1,2-Dibromoethane	0.586	0.507	13.5	86	0.00
55 I	Chlorobenzene-d5	1.000	1.000	0.0	88	0.00
56	1,1,1,2-Tetrachloroethane	0.605	0.573	5.3	92	0.00
57 T	Chlorobenzene	1.040	0.930	10.6	86	0.00
58 T	Ethyl Benzene	1.981	1.743	12.0	84	0.00
59 T	m/p-Xylene	1.596	1.426	10.7	86	0.00
60 T	o-Xylene	1.607	1.459	9.2	89	0.00
61 T	Styrene	0.767	0.649	15.4	79	0.00
62	Isopropylbenzene	2.360	2.105	10.8	85	0.00
63 T	1,1,2,2-Tetrachloroethane	1.042	0.842	19.2	85	0.00
64	n-propylbenzene	0.593	0.519	12.5	84	0.00
65	tert-Butylbenzene	2.169	1.920	11.5	88	0.00
66 T	Benzyl Chloride	0.238	0.192	19.3	69	0.00
67	sec-Butylbenzene	3.017	2.685	11.0	89	0.00
68 S	1-Bromo-4-Fluorobenzene	0.822	0.882	-7.3	93	0.00
69	p-Isopropyltoluene	2.621	2.303	12.1	88	0.00
70	n-Butylbenzene	2.730	2.423	11.2	90	0.00
71	2-Chlorotoluene	1.851	1.660	10.3	86	0.00
72 T	4-Ethyltoluene	1.965	1.755	10.7	87	0.00
73 T	1,3,5-Trimethylbenzene	1.735	1.518	12.5	88	0.00
74 T	1,2,4-Trimethylbenzene	2.026	1.690	16.6	89	0.00
75 T	1,3-Dichlorobenzene	1.053	0.888	15.7	82	0.00
76 T	1,4-Dichlorobenzene	1.059	0.894	15.6	83	0.00
77 T	1,2-Dichlorobenzene	1.024	0.875	14.6	85	0.00
78 T	Hexachloro-1,3-Butadiene	1.034	0.828	19.9	84	0.00
79 T	Naphthalene	2.423	1.920	20.8	84	0.00
80 T	Naphthalene,2-methyl-	1.381	1.023	25.9	81	0.00
81 T	1,2,4-Trichlorobenzene	1.040	0.844	18.8	82	0.00

(#) = Out of Range

SPCC's out = 0 CCC's out = 0

Data Path : Z:\voasrv\HPCHEM1\MSVOA_L\Data\VL042125\
 Data File : VL042395.D
 Acq On : 21 Apr 2025 08:41
 Operator : SY/MD
 Sample : VSTDCCC010
 Misc : 400mL/MSVOA_L
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_L
 LabSampleId :
 VSTDCCC010

Quant Time: Apr 22 02:50:13 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_L\methods\VL041725AIR.M
 Quant Title : AIR ANALYSIS BY METHOD TO-15 Instrument: MSVOA_L Fri Aug 26 06:05:16 2022
 QLast Update : Fri Apr 18 02:14:00 2025
 Response via : Initial Calibration

Min. RRF : 0.050 Min. Rel. Area : 20% Max. R.T. Dev 0.50min
 Max. RRF Dev : 30% Max. Rel. Area : 150%

	Compound	Amount	Calc.	%Dev	Area%	Dev(min)
1 I	Bromochloromethane	10.000	10.000	0.0	94	0.00
2 T	Dichlorodifluoromethane	10.000	10.253	-2.5	103	0.00
3	Chlorodifluoromethane	10.000	8.652	13.5	87	0.00
4	Chloromethane	10.000	9.455	5.4	96	0.00
5 T	Vinyl Chloride	10.000	8.736	12.6	98	0.00
6 T	Bromomethane	10.000	9.484	5.2	96	0.00
7	Chloroethane	10.000	9.806	1.9	101	0.00
8 T	Dichlorotetrafluoroethane	10.000	10.155	-1.5	101	0.00
9 T	Propene	10.000	7.422	25.8	78	0.00
10 T	Heptane	10.000	7.026	29.7	76	0.00
11 T	Trichlorofluoromethane	10.000	10.328	-3.3	104	0.00
12 T	1,1,2-Trichlorotrifluoroeth	10.000	9.515	4.8	96	0.00
13	Ethanol	10.000	3.130	68.7#	102	0.00
14 T	Bromoethene	10.000	9.680	3.2	96	0.00
15 T	Acetone	10.000	8.675	13.2	106	0.00
16 T	1,3-Butadiene	10.000	8.762	12.4	99	0.00
17	tert-Butyl alcohol	10.000	9.218	7.8	98	0.00
18 T	1,1-Dichloroethene	10.000	9.244	7.6	96	0.00
19 T	Isopropyl Alcohol	10.000	9.126	8.7	95	0.00
20 T	Methylene Chloride	10.000	7.915	20.8	96	0.00
21 T	Allyl Chloride	10.000	9.056	9.4	99	0.00
22 T	trans-1,2-Dichloroethene	10.000	9.192	8.1	95	0.00
23 T	Vinyl Acetate	10.000	7.306	26.9	82	0.00
24 T	1,1-Dichloroethane	10.000	9.405	6.0	96	0.00
25 T	Ethyl Acetate	10.000	8.211	17.9	85	0.00
26 T	Hexane	10.000	7.789	22.1	84	0.00
27 T	Carbon Disulfide	10.000	9.658	3.4	96	0.00
28 T	Methyl tert-Butyl Ether	10.000	8.414	15.9	90	0.00
29 T	Chloroform	10.000	9.502	5.0	96	0.00
30 T	Cyclohexane	10.000	8.095	19.0	81	0.00
31 T	cis-1,2-Dichloroethene	10.000	9.051	9.5	91	0.00
32 T	1,1,1-Trichloroethane	10.000	9.284	7.2	100	0.00
33 I	1,4-Difluorobenzene	10.000	10.000	0.0	91	0.00
34 T	2-Butanone	10.000	8.316	16.8	87	0.00
35 T	Carbon Tetrachloride	10.000	9.614	3.9	103	0.00
36 T	Benzene	10.000	8.420	15.8	83	0.00
37 T	1,2-Dichloroethane	10.000	10.206	-2.1	102	0.00
38 T	Trichloroethene	10.000	7.804	22.0	85	0.00
39 T	1,2-Dichloropropane	10.000	8.018	19.8	80	0.00
40 T	1,4-Dioxane	10.000	7.286	27.1	81	0.00
41 T	Tetrahydrofuran	10.000	7.701	23.0	79	0.00
42 T	Bromodichloromethane	10.000	9.800	2.0	95	0.00
43	Methyl Methacrylate	10.000	7.667	23.3	79	0.00
44 T	2,2,4-Trimethylpentane	10.000	8.138	18.6	81	-0.01
45 T	t-1,3-Dichloropropene	10.000	8.596	14.0	85	0.00
46 T	cis-1,3-Dichloropropene	10.000	8.595	14.0	84	0.00
47 T	1,1,2-Trichloroethane	10.000	8.711	12.9	87	0.00
48 T	Dibromochloromethane	10.000	9.476	5.2	92	0.00

Data Path : Z:\voasrv\HPCHEM1\MSVOA_L\Data\VL042125\
 Data File : VL042395.D
 Acq On : 21 Apr 2025 08:41
 Operator : SY/MD
 Sample : VSTDCCC010
 Misc : 400mL/MSVOA_L
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_L
 LabSampleId :
 VSTDCCC010

Quant Time: Apr 22 02:50:13 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_L\methods\VL041725AIR.M
 Quant Title : AIR ANALYSIS BY METHOD TO-15 Instrument: MSVOA_L Fri Aug 26 06:05:16 2022
 QLast Update : Fri Apr 18 02:14:00 2025
 Response via : Initial Calibration

Min. RRF : 0.050 Min. Rel. Area : 20% Max. R.T. Dev 0.50min
 Max. RRF Dev : 30% Max. Rel. Area : 150%

	Compound	Amount	Calc.	%Dev	Area%	Dev(min)
49 T	Bromoform	10.000	8.836	11.6	83	0.00
50 T	4-Methyl-2-Pentanone	10.000	7.314	26.9	81	0.00
51 T	2-Hexanone	10.000	6.590	34.1#	79	0.00
52 T	Tetrachloroethene	10.000	7.474	25.3	80	0.00
53 T	Toluene	10.000	8.206	17.9	81	0.00
54 T	1,2-Dibromoethane	10.000	8.654	13.5	86	0.00
55 I	Chlorobenzene-d5	10.000	10.000	0.0	88	0.00
56	1,1,1,2-Tetrachloroethane	10.000	9.468	5.3	92	0.00
57 T	Chlorobenzene	10.000	8.944	10.6	86	0.00
58 T	Ethyl Benzene	10.000	8.797	12.0	84	0.00
59 T	m/p-Xylene	20.000	17.868	10.7	86	0.00
60 T	o-Xylene	10.000	9.078	9.2	89	0.00
61 T	Styrene	10.000	8.454	15.5	79	0.00
62	Isopropylbenzene	10.000	8.920	10.8	85	0.00
63 T	1,1,2,2-Tetrachloroethane	10.000	8.085	19.1	85	0.00
64	n-propylbenzene	10.000	8.744	12.6	84	0.00
65	tert-Butylbenzene	10.000	8.853	11.5	88	0.00
66 T	Benzyl Chloride	10.000	8.061	19.4	69	0.00
67	sec-Butylbenzene	10.000	8.899	11.0	89	0.00
68 S	1-Bromo-4-Fluorobenzene	10.000	10.736	-7.4	93	0.00
69	p-Isopropyltoluene	10.000	8.789	12.1	88	0.00
70	n-Butylbenzene	10.000	8.877	11.2	90	0.00
71	2-Chlorotoluene	10.000	8.967	10.3	86	0.00
72 T	4-Ethyltoluene	10.000	8.929	10.7	87	0.00
73 T	1,3,5-Trimethylbenzene	10.000	8.752	12.5	88	0.00
74 T	1,2,4-Trimethylbenzene	10.000	8.340	16.6	89	0.00
75 T	1,3-Dichlorobenzene	10.000	8.436	15.6	82	0.00
76 T	1,4-Dichlorobenzene	10.000	8.445	15.5	83	0.00
77 T	1,2-Dichlorobenzene	10.000	8.548	14.5	85	0.00
78 T	Hexachloro-1,3-Butadiene	10.000	8.002	20.0	84	0.00
79 T	Naphthalene	10.000	7.925	20.8	84	0.00
80 T	Naphthalene,2-methyl-	10.000	7.412	25.9	81	0.00
81 T	1,2,4-Trichlorobenzene	10.000	8.112	18.9	82	0.00

(#) = Out of Range

SPCC's out = 0 CCC's out = 0



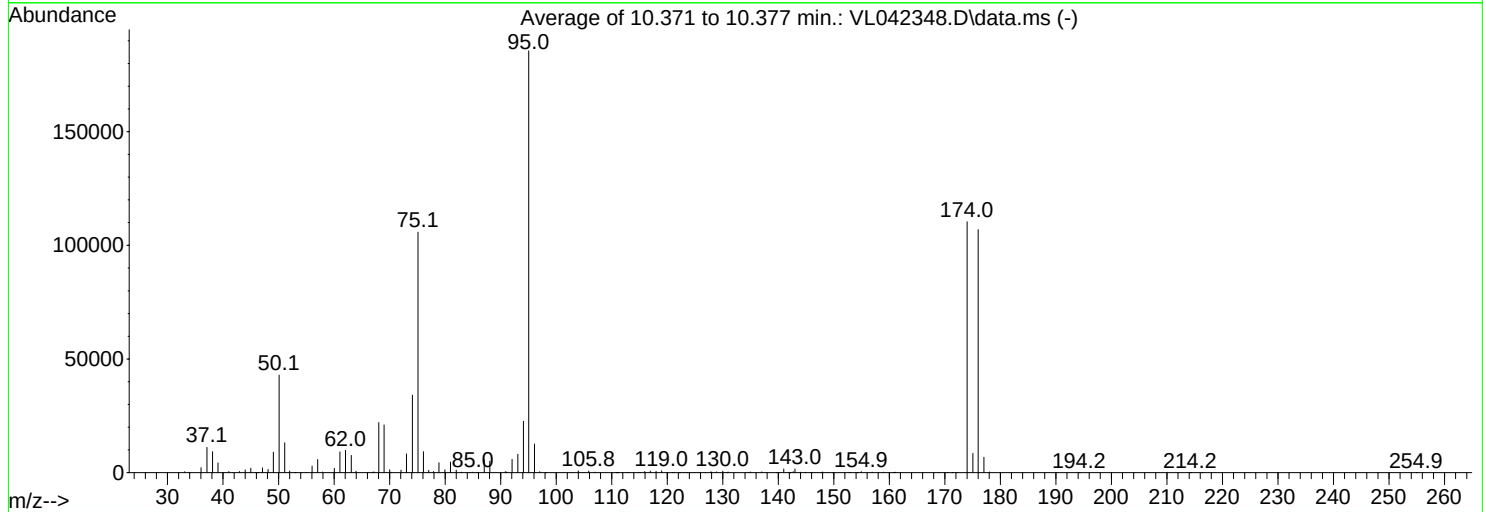
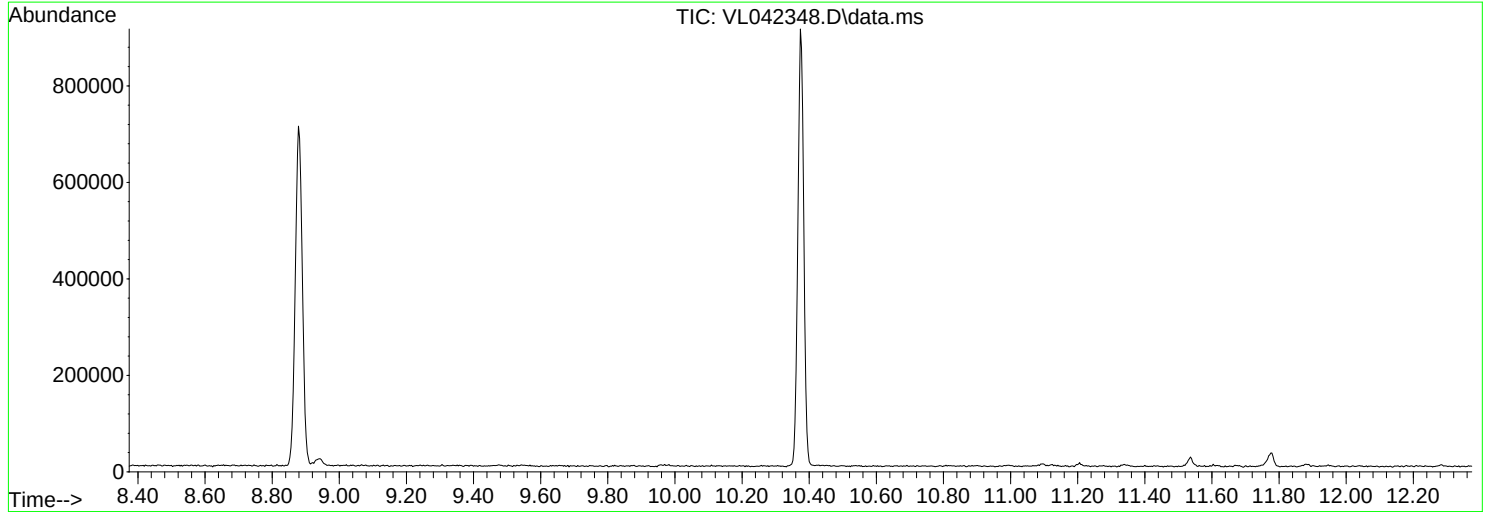
QC SAMPLE DATA

Data Path : Z:\voasrv\HPCHEM1\MSVOA_L\Data\VL041725\
Data File : VL042348.D
Acq On : 17 Apr 2025 07:50
Operator : SY/MD
Sample : BFB
Misc : 400mL/MSVOA_L
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_L
ClientSampleId :
BFB

Integration File: RTEINT.P

Method : Z:\voasrv\HPCHEM1\MSVOA_L\methods\VL041725AIR.M
Title : AIR ANALYSIS BY METHOD TO-15 Instrument: MSVOA_L Fri Aug 26 06:05:16 2022
Last Update : Fri Apr 18 02:14:00 2025



AutoFind: Scans 3177, 3178, 3179; Background Corrected with Scan 3162

Target Mass	Rel. to Mass	Lower Limit%	Upper Limit%	Rel. Abn%	Raw Abn	Result Pass/Fail
50	95	8	40	23.1	42933	PASS
75	95	30	66	57.0	105781	PASS
95	95	100	100	100.0	185564	PASS
96	95	5	9	6.8	12627	PASS
173	174	0.00	2	0.0	0	PASS
174	95	50	120	59.5	110360	PASS
175	174	4	9	7.6	8420	PASS
176	174	93	101	96.9	106933	PASS
177	176	5	9	6.3	6744	PASS

Average of 10.371 to 10.377 min.: VL042348.D\data.ms

BFB

Modified:subtracted

m/z	abund.	m/z	abund.	m/z	abund.	m/z	abund.
33.10	400	45.00	1996	55.00	78	67.20	32
34.00	54	45.90	94	56.05	2900	68.05	2203
36.05	2214	46.20	34	57.05	5754	69.00	2101
37.10	11145	47.10	2100	57.95	265	70.00	130
38.10	9197	48.10	1438	60.05	1919	70.20	209
39.10	4289	49.05	9045	61.05	9155	72.05	1141
39.95	158	50.10	42933	62.05	9788	73.05	8240
41.05	513	51.10	13188	63.10	7593	74.10	34064
41.80	79	52.00	718	63.95	632	75.10	105781
42.95	504	52.95	9	66.10	115	76.10	9240
44.00	1207	54.15	112	66.90	196	77.05	979

Average of 10.371 to 10.377 min.: VL042348.D\data.ms

BFB

Modified:subtracted

m/z	abund.	m/z	abund.	m/z	abund.	m/z	abund.
77.90	494	90.85	447	105.00	95	115.95	528
78.90	4344	91.10	262	105.85	680	116.95	712
79.95	1296	92.05	5876	107.00	225	118.00	575
80.95	4648	93.10	8069	110.10	136	119.00	881
82.00	1107	94.10	22619	110.80	37	120.10	70
83.00	249	95.05	185564	111.10	8	120.90	30
84.40	55	96.10	12627	111.95	156	121.30	47
85.00	227	97.05	457	112.75	77	122.20	21
85.85	131	98.05	5	113.10	144	122.95	44
87.05	5767	102.90	78	114.80	135	123.50	25
88.05	5057	104.00	767	115.10	50	124.10	89

Average of 10.371 to 10.377 min.: VL042348.D\data.ms

BFB

Modified:subtracted

m/z	abund.	m/z	abund.	m/z	abund.	m/z	abund.
124.90	71	137.05	364	146.15	158	153.90	93
125.75	72	139.65	74	146.65	58	154.20	27
126.90	80	140.00	38	147.10	68	154.95	399
127.80	297	141.00	1612	147.70	38	156.90	228
128.00	275	141.80	139	148.00	288	158.95	203
128.85	366	142.10	34	148.95	115	160.80	26
130.00	693	142.70	434	149.80	37	161.05	186
130.80	55	143.00	1656	149.95	79	168.90	83
130.95	285	143.95	200	151.90	60	170.50	119
134.70	94	144.95	161	152.75	91	171.30	19
135.00	292	145.80	67	153.05	144	171.90	90

Average of 10.371 to 10.377 min.: VL042348.D\data.ms

BFB

Modified:subtracted

m/z	abund.	m/z	abund.	m/z	abund.	m/z	abund.
172.30	132						
174.00	110360						
175.05	8420						
176.00	106933						
177.05	6744						
178.05	298						
194.20	17						
214.20	17						
216.30	17						
254.90	17						

Instrument :

MSVOA_L

ClientSampleId :

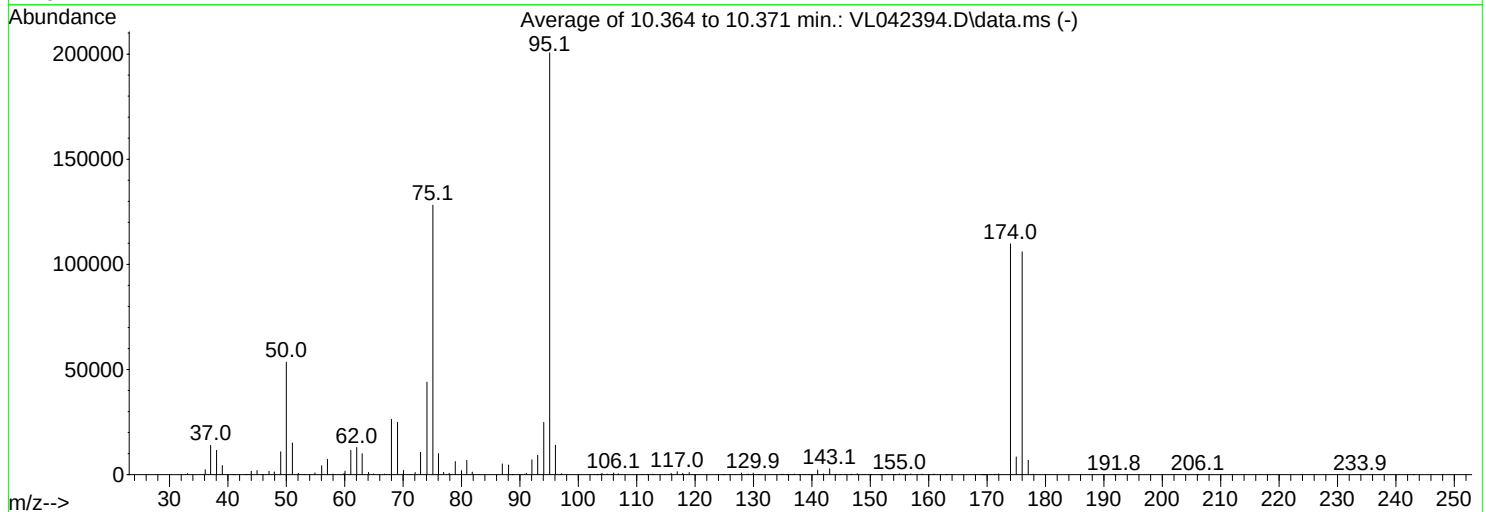
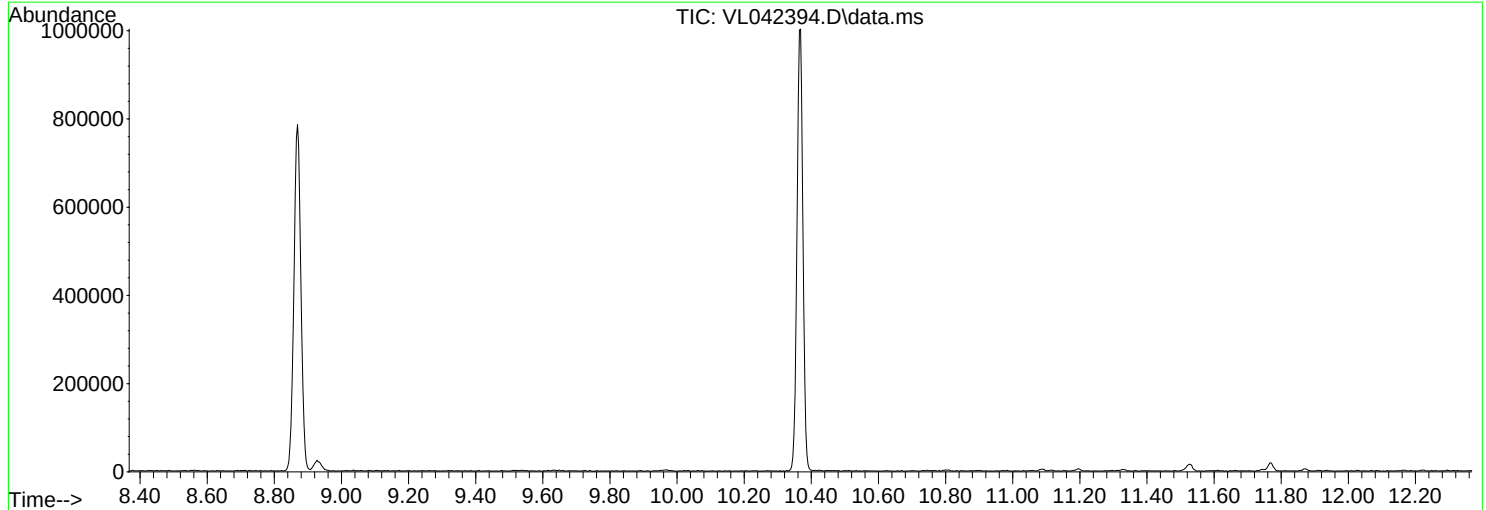
BFB

Data Path : Z:\voasrv\HPCHEM1\MSVOA_L\Data\VL042125\
Data File : VL042394.D
Acq On : 21 Apr 2025 07:58
Operator : SY/MD
Sample : BFB
Misc : 400mL/MSVOA_L
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_L
ClientSampleId :
BFB

Integration File: RTEINT.P

Method : Z:\voasrv\HPCHEM1\MSVOA_L\methods\VL041725AIR.M
Title : AIR ANALYSIS BY METHOD TO-15 Instrument: MSVOA_L Fri Aug 26 06:05:16 2022
Last Update : Fri Apr 18 02:14:00 2025



AutoFind: Scans 3175, 3176, 3177; Background Corrected with Scan 3159

Target	Rel. to	Lower	Upper	Rel.	Raw	Result
Mass	Mass	Limit%	Limit%	Abn%	Abn	Pass/Fail
50	95	8	40	26.7	53502	PASS
75	95	30	66	63.9	128181	PASS
95	95	100	100	100.0	200597	PASS
96	95	5	9	7.0	13990	PASS
173	174	0.00	2	0.0	0	PASS
174	95	50	120	54.7	109757	PASS
175	174	4	9	7.7	8464	PASS
176	174	93	101	96.6	106032	PASS
177	176	5	9	6.4	6826	PASS

Average of 10.364 to 10.371 min.: VL042394.D\data.ms

BFB

Modified:subtracted

m/z	abund.	m/z	abund.	m/z	abund.	m/z	abund.
33.10	518	45.00	2010	54.90	832	66.85	304
36.10	2394	45.85	194	56.05	4237	67.20	161
37.05	13953	46.10	197	57.05	7329	68.00	2631
38.05	11556	47.05	1570	58.00	328	69.05	2485
39.05	4320	47.95	1283	59.90	593	70.05	2000
39.90	57	49.05	10899	60.05	1702	72.05	1049
40.90	31	50.00	53502	61.05	11547	73.00	10643
41.30	48	51.05	15046	62.05	12934	74.10	44016
43.00	64	52.05	608	63.00	9941	75.10	128181
43.20	184	52.70	76	64.10	1108	76.05	10068
44.00	1632	53.15	138	64.90	355	76.95	1033

Average of 10.364 to 10.371 min.: VL042394.D\data.ms

BFB

Modified:subtracted

m/z	abund.	m/z	abund.	m/z	abund.	m/z	abund.
77.90	700	91.10	697	104.05	598	114.80	178
78.95	6226	92.05	7067	104.95	278	115.90	632
80.00	1878	93.05	9194	105.70	188	116.95	1450
80.90	6765	94.10	24813	106.05	831	117.90	579
81.85	1186	95.10	200597	106.85	467	119.00	1019
83.00	184	96.10	13990	109.95	82	121.90	17
86.05	106	96.90	128	110.85	113	124.00	56
87.00	5101	97.15	433	111.10	91	124.80	25
88.05	4600	102.75	176	111.85	167	125.90	42
88.70	40	103.20	26	112.75	225	126.85	64
90.80	235	103.80	263	113.00	54	127.95	770

Average of 10.364 to 10.371 min.: VL042394.D\data.ms

BFB

Modified:subtracted

m/z	abund.	m/z	abund.	m/z	abund.	m/z	abund.
128.80	183	137.80	21	145.00	70	152.60	135
129.10	275	138.70	105	145.70	127	153.10	56
129.95	780	138.90	54	145.85	120	153.80	154
130.85	125	139.55	52	146.75	211	154.20	31
131.00	55	140.05	163	147.75	498	154.80	81
131.20	65	141.00	2199	148.80	55	155.00	397
134.80	172	141.95	402	149.05	129	155.70	99
135.05	261	143.05	2809	149.65	112	156.20	25
135.70	20	144.00	151	150.10	20	156.95	371
136.60	73	144.20	61	151.70	27	158.85	322
137.00	333	144.80	162	152.00	33	160.90	347

Average of 10.364 to 10.371 min.: VL042394.D\data.ms

BFB

Modified:subtracted

m/z	abund.	m/z	abund.	m/z	abund.	m/z	abund.
163.00	37	176.00	106032				
169.00	59	177.05	6826				
169.20	38	177.90	242				
169.50	29	191.75	41				
169.70	25	206.10	24				
170.65	116	233.90	67				
171.20	28	243.00	24				
171.45	164						
171.95	381						
174.00	109757						
175.00	8464						

Report of Analysis

Client:	GFE LLC	Date Collected:	
Project:	416 Clinton St Hempstead	Date Received:	
Client Sample ID:	VL0421ABL01	SDG No.:	Q1801
Lab Sample ID:	VL0421ABL01	Matrix:	Air
Analytical Method:	TO-15	Test:	TO-15
Sample Wt/Vol:	400	Units:	mL

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
VL042396.D	1		04/21/25 09:28	VL042125

CAS Number	Parameter	Conc. ppbv	Conc. ug/M3	Qualifier	MDL	LOQ / CRQL	Units
TARGETS							
75-71-8	Dichlorodifluoromethane	0.21	1.04	U	1.04	2.47	ug/m3
74-87-3	Chloromethane	0.10	0.21	U	0.21	1.03	ug/m3
75-01-4	Vinyl Chloride	0.010	0.030	U	0.030	0.080	ug/m3
74-83-9	Bromomethane	0.14	0.54	U	0.54	1.94	ug/m3
75-00-3	Chloroethane	0.17	0.45	U	0.45	1.32	ug/m3
109-99-9	Tetrahydrofuran	0.13	0.38	U	0.38	1.47	ug/m3
75-69-4	Trichlorofluoromethane	0.16	0.90	U	0.90	2.81	ug/m3
76-13-1	1,1,2-Trichlorotrifluoroethane	0.16	1.23	U	1.23	3.83	ug/m3
76-14-2	Dichlorotetrafluoroethane	0.090	0.63	U	0.63	3.49	ug/m3
75-65-0	tert-Butyl alcohol	0.19	0.58	U	0.58	1.52	ug/m3
142-82-5	Heptane	0.11	0.45	U	0.45	2.05	ug/m3
75-35-4	1,1-Dichloroethene	0.14	0.56	U	0.56	1.98	ug/m3
67-64-1	Acetone	0.24	0.57	U	0.57	1.19	ug/m3
75-15-0	Carbon Disulfide	0.16	0.50	U	0.50	1.56	ug/m3
1634-04-4	Methyl tert-Butyl Ether	0.12	0.43	U	0.43	1.80	ug/m3
75-09-2	Methylene Chloride	0.19	0.66	U	0.66	1.74	ug/m3
156-60-5	trans-1,2-Dichloroethene	0.19	0.75	U	0.75	1.98	ug/m3
75-34-3	1,1-Dichloroethane	0.13	0.53	U	0.53	2.02	ug/m3
110-82-7	Cyclohexane	0.22	0.76	U	0.76	1.72	ug/m3
78-93-3	2-Butanone	0.12	0.35	U	0.35	1.47	ug/m3
56-23-5	Carbon Tetrachloride	0.010	0.060	U	0.060	0.19	ug/m3
156-59-2	cis-1,2-Dichloroethene	0.090	0.36	U	0.36	1.98	ug/m3
67-66-3	Chloroform	0.080	0.39	U	0.39	2.44	ug/m3
71-55-6	1,1,1-Trichloroethane	0.010	0.050	U	0.050	0.16	ug/m3
540-84-1	2,2,4-Trimethylpentane	0.10	0.47	U	0.47	2.34	ug/m3
71-43-2	Benzene	0.090	0.29	U	0.29	1.60	ug/m3
107-06-2	1,2-Dichloroethane	0.090	0.36	U	0.36	2.02	ug/m3
79-01-6	Trichloroethene	0.020	0.11	U	0.11	0.16	ug/m3
78-87-5	1,2-Dichloropropane	0.11	0.51	U	0.51	2.31	ug/m3
75-27-4	Bromodichloromethane	0.080	0.54	U	0.54	3.35	ug/m3
108-10-1	4-Methyl-2-Pentanone	0.090	0.37	U	0.37	2.05	ug/m3
108-88-3	Toluene	0.11	0.41	U	0.41	1.88	ug/m3
10061-02-6	t-1,3-Dichloropropene	0.060	0.27	U	0.27	2.27	ug/m3
10061-01-5	cis-1,3-Dichloropropene	0.060	0.27	U	0.27	2.27	ug/m3
79-00-5	1,1,2-Trichloroethane	0.070	0.38	U	0.38	2.73	ug/m3

Report of Analysis

Client:	GFE LLC	Date Collected:	
Project:	416 Clinton St Hempstead	Date Received:	
Client Sample ID:	VL0421ABL01	SDG No.:	Q1801
Lab Sample ID:	VL0421ABL01	Matrix:	Air
Analytical Method:	TO-15	Test:	TO-15
Sample Wt/Vol:	400	Units:	mL

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
VL042396.D	1		04/21/25 09:28	VL042125

CAS Number	Parameter	Conc. ppbv	Conc. ug/M3	Qualifier	MDL	LOQ / CRQL	Units
124-48-1	Dibromochloromethane	0.070	0.60	U	0.60	4.26	ug/m3
106-93-4	1,2-Dibromoethane	0.070	0.54	U	0.54	0.77	ug/m3
127-18-4	Tetrachloroethene	0.020	0.14	U	0.14	0.20	ug/m3
108-90-7	Chlorobenzene	0.080	0.37	U	0.37	2.30	ug/m3
100-41-4	Ethyl Benzene	0.12	0.52	U	0.52	2.17	ug/m3
179601-23-1	m/p-Xylene	0.21	0.91	U	0.91	4.34	ug/m3
95-47-6	o-Xylene	0.12	0.52	U	0.52	2.17	ug/m3
100-42-5	Styrene	0.11	0.47	U	0.47	2.13	ug/m3
75-25-2	Bromoform	0.060	0.62	U	0.62	5.17	ug/m3
79-34-5	1,1,2,2-Tetrachloroethane	0.010	0.070	U	0.070	0.21	ug/m3
95-49-8	2-Chlorotoluene	0.11	0.57	U	0.57	2.59	ug/m3
108-67-8	1,3,5-Trimethylbenzene	0.11	0.54	U	0.54	2.46	ug/m3
95-63-6	1,2,4-Trimethylbenzene	0.080	0.39	U	0.39	2.46	ug/m3
541-73-1	1,3-Dichlorobenzene	0.080	0.48	U	0.48	3.01	ug/m3
106-46-7	1,4-Dichlorobenzene	0.050	0.30	U	0.30	3.01	ug/m3
95-50-1	1,2-Dichlorobenzene	0.080	0.48	U	0.48	3.01	ug/m3
120-82-1	1,2,4-Trichlorobenzene	0.090	0.67	U	0.67	3.71	ug/m3
87-68-3	Hexachloro-1,3-Butadiene	0.090	0.96	U	0.96	5.33	ug/m3
106-99-0	1,3-Butadiene	0.13	0.29	U	0.29	1.11	ug/m3
91-20-3	Naphthalene	0.080	0.42	U	0.42	0.52	ug/m3
622-96-8	4-Ethyltoluene	0.12	0.59	U	0.59	2.46	ug/m3
110-54-3	Hexane	0.11	0.39	U	0.39	1.76	ug/m3
107-05-1	Allyl Chloride	0.15	0.47	U	0.47	1.57	ug/m3
123-91-1	1,4-Dioxane	0.21	0.76	U	0.76	1.80	ug/m3
80-62-6	Methyl Methacrylate	0.090	0.37	U	0.37	2.05	ug/m3
SURROGATES							
460-00-4	1-Bromo-4-Fluorobenzene	10.5			65 - 135	105%	SPK: 10
INTERNAL STANDARDS							
74-97-5	Bromochloromethane	120000		2.784			
540-36-3	1,4-Difluorobenzene	342000		3.952			
3114-55-4	Chlorobenzene-d5	294000		8.878			

Report of Analysis

Client:	GFE LLC	Date Collected:	
Project:	416 Clinton St Hempstead	Date Received:	
Client Sample ID:	VL0421ABL01	SDG No.:	Q1801
Lab Sample ID:	VL0421ABL01	Matrix:	Air
Analytical Method:	TO-15	Test:	TO-15
Sample Wt/Vol:	400	Units:	mL

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
VL042396.D	1		04/21/25 09:28	VL042125

CAS Number	Parameter	Conc. ppbv	Conc. ug/M3	Qualifier	MDL	LOQ / CRQL	Units
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U = Not Detected
 RL = Reporting Limit
 MDL = Method Detection Limit
 E = Value Exceeds Calibration Range
 D = Dilution

J = Estimated Value
 B = Analyte Found in Associated Method Blank
 N = Presumptive Evidence of a Compound
 * = Values outside of QC limits
 Q = indicates LCS control criteria did not meet requirements

Data Path : Z:\voasrv\HPCHEM1\MSVOA_L\Data\VL042125\
 Data File : VL042396.D
 Acq On : 21 Apr 2025 09:28
 Operator : SY/MD
 Sample : VL0421ABL01
 Misc : 400mL/MSVOA_L
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_L
 ClientSampleId :
 VL0421ABL01

Quant Time: Apr 22 02:50:57 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_L\methods\VL041725AIR.M
 Quant Title : AIR ANALYSIS BY METHOD TO-15 Instrument: MSVOA_L Fri Aug 26 06:05:16 2022
 QLast Update : Fri Apr 18 02:14:00 2025
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) Bromochloromethane	2.784	49	120411	10.000	ppbv	0.00
33) 1,4-Difluorobenzene	3.952	114	341628	10.000	ppbv	0.00
55) Chlorobenzene-d5	8.878	117	293743	10.000	ppbv	0.00

System Monitoring Compounds

68) 1-Bromo-4-Fluorobenzene	10.374	95	252683	10.468	ppbv	0.00
Spiked Amount	10.000	Range	65 - 135	Recovery	=	104.700%

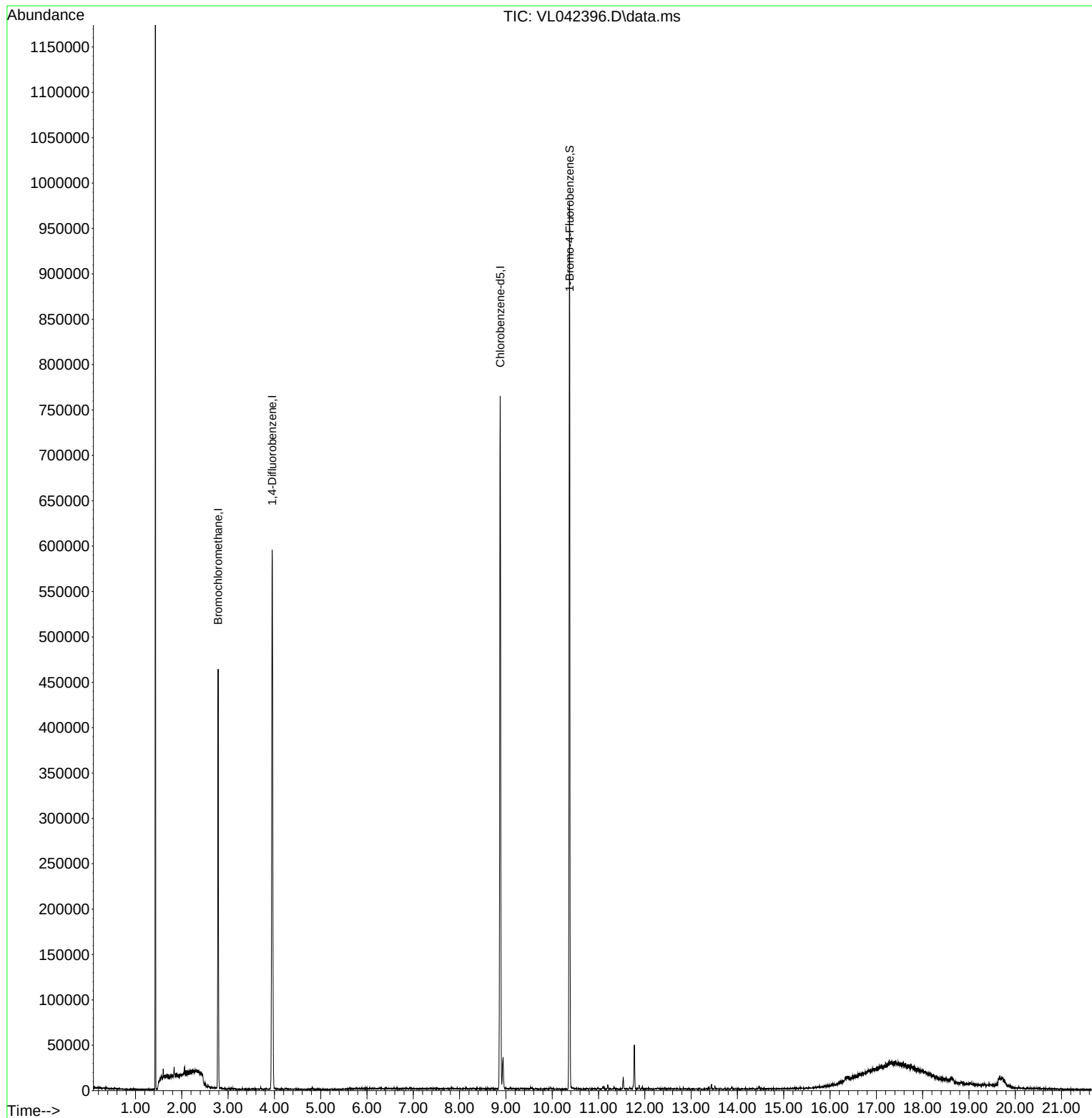
Target Compounds	Qvalue

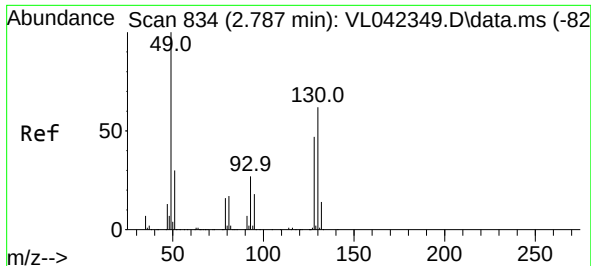
(#) = qualifier out of range (m) = manual integration (+) = signals summed

Data Path : Z:\voasrv\HPCHEM1\MSVOA_L\Data\VL042125\
Data File : VL042396.D
Acq On : 21 Apr 2025 09:28
Operator : SY/MD
Sample : VL0421ABL01
Misc : 400mL/MSVOA_L
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_L
ClientSampleId :
VL0421ABL01

Quant Time: Apr 22 02:50:57 2025
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_L\methods\VL041725AIR.M
Quant Title : AIR ANALYSIS BY METHOD TO-15 Instrument: MSVOA_L Fri Aug 26 06:05:16 2022
QLast Update : Fri Apr 18 02:14:00 2025
Response via : Initial Calibration

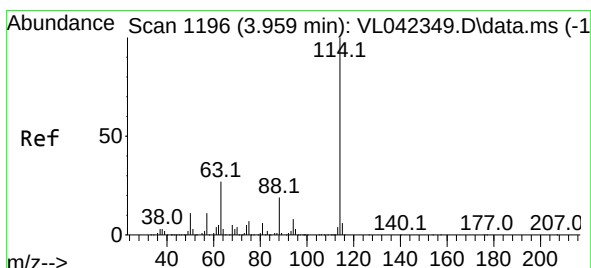
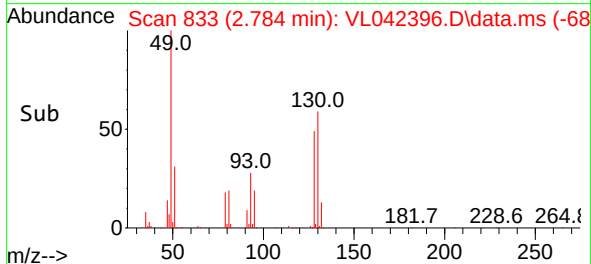
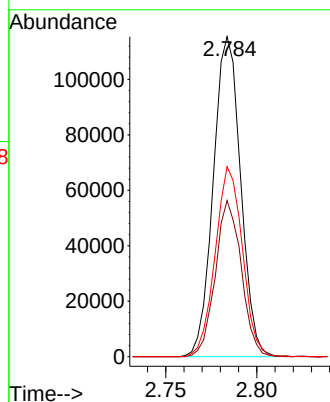
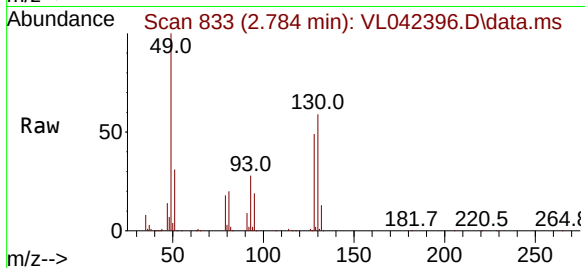




#1
Bromochloromethane
Concen: 10.000 ppbv
RT: 2.784 min Scan# 81
Delta R.T. -0.003 min
Lab File: VL042396.D
Acq: 21 Apr 2025 09:28

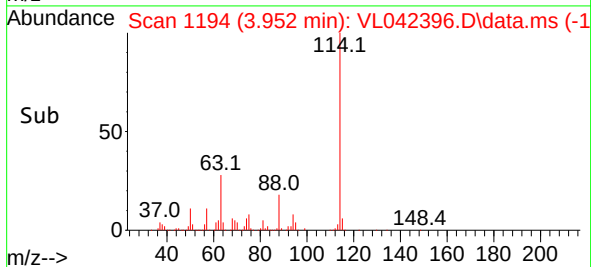
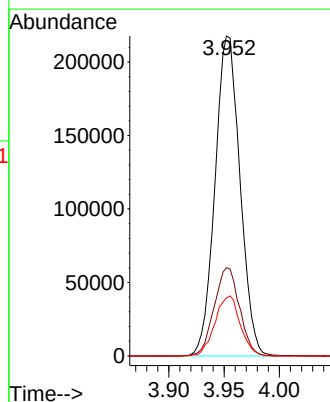
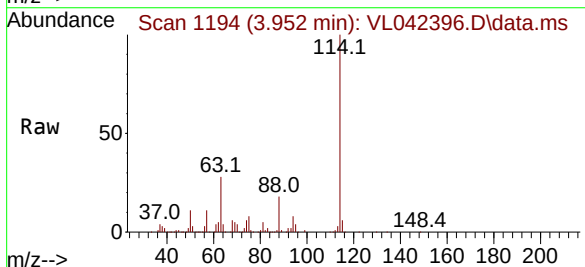
Instrument :
MSVOA_L
ClientSampleId :
VL0421ABL01

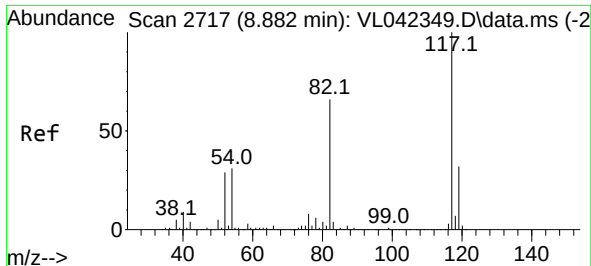
Tgt Ion: 49 Resp: 120411
Ion Ratio Lower Upper
49 100
128 46.4 22.8 68.4
130 59.3 29.3 88.0



#33
1,4-Difluorobenzene
Concen: 10.000 ppbv
RT: 3.952 min Scan# 1194
Delta R.T. -0.007 min
Lab File: VL042396.D
Acq: 21 Apr 2025 09:28

Tgt Ion: 114 Resp: 341628
Ion Ratio Lower Upper
114 100
63 27.9 21.8 32.6
88 19.0 15.5 23.3

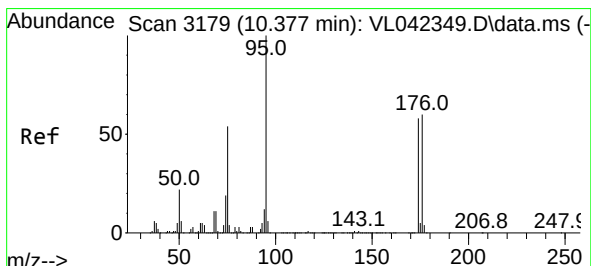
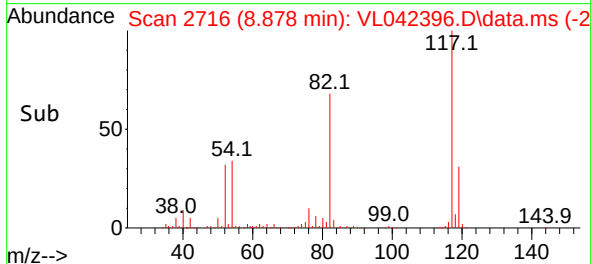
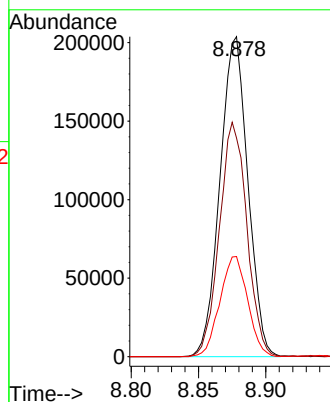
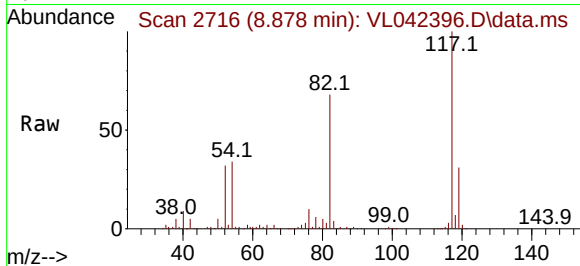




#55
Chlorobenzene-d5
Concen: 10.000 ppbv
RT: 8.878 min Scan# 21
Delta R.T. -0.003 min
Lab File: VL042396.D
Acq: 21 Apr 2025 09:28

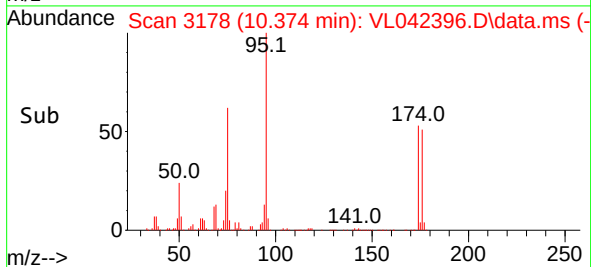
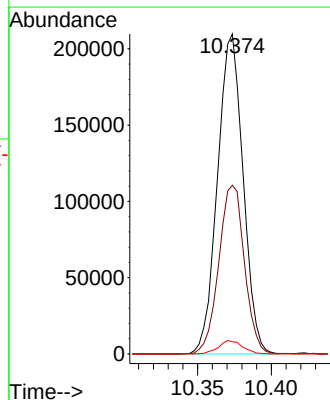
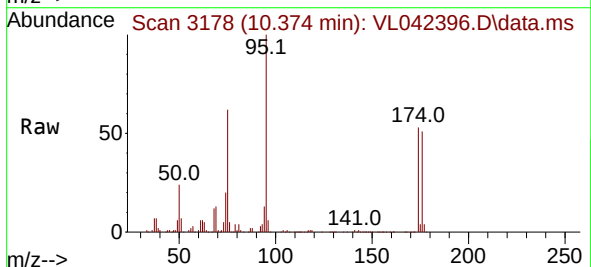
Instrument :
MSVOA_L
ClientSampleId :
VL0421ABL01

Tgt Ion:117 Resp: 293743
Ion Ratio Lower Upper
117 100
82 72.2 57.8 86.6
119 31.3 25.2 37.8



#68
1-Bromo-4-Fluorobenzene
Concen: 10.468 ppbv
RT: 10.374 min Scan# 3178
Delta R.T. -0.003 min
Lab File: VL042396.D
Acq: 21 Apr 2025 09:28

Tgt Ion: 95 Resp: 252683
Ion Ratio Lower Upper
95 100
174 53.1 48.5 72.7
175 4.0 3.7 5.5



Report of Analysis

Client:	GFE LLC	Date Collected:	
Project:	416 Clinton St Hempstead	Date Received:	
Client Sample ID:	VL0421ABS01	SDG No.:	Q1801
Lab Sample ID:	VL0421ABS01	Matrix:	Air
Analytical Method:	TO-15	Test:	TO-15
Sample Wt/Vol:	400	Units:	mL

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
VL042397.D	1		04/21/25 10:14	VL042125

CAS Number	Parameter	Conc. ppbv	Conc. ug/M3	Qualifier	MDL	LOQ / CRQL	Units
TARGETS							
75-71-8	Dichlorodifluoromethane	10.2	50.4		1.04	2.47	ug/m3
74-87-3	Chloromethane	9.60	19.8		0.21	1.03	ug/m3
75-01-4	Vinyl Chloride	8.60	22.0		0.030	0.080	ug/m3
74-83-9	Bromomethane	9.10	35.3		0.54	1.94	ug/m3
75-00-3	Chloroethane	9.50	25.1		0.45	1.32	ug/m3
109-99-9	Tetrahydrofuran	7.80	23.0		0.38	1.47	ug/m3
75-69-4	Trichlorofluoromethane	10.3	57.9		0.90	2.81	ug/m3
76-13-1	1,1,2-Trichlorotrifluoroethane	9.60	73.6		1.23	3.83	ug/m3
76-14-2	Dichlorotetrafluoroethane	10.1	70.6		0.63	3.49	ug/m3
75-65-0	tert-Butyl alcohol	9.00	27.3		0.58	1.52	ug/m3
142-82-5	Heptane	7.30	29.9		0.45	2.05	ug/m3
75-35-4	1,1-Dichloroethene	9.40	37.3		0.56	1.98	ug/m3
67-64-1	Acetone	8.70	20.7		0.57	1.19	ug/m3
75-15-0	Carbon Disulfide	9.40	29.3		0.50	1.56	ug/m3
1634-04-4	Methyl tert-Butyl Ether	8.50	30.6		0.43	1.80	ug/m3
75-09-2	Methylene Chloride	8.00	27.8		0.66	1.74	ug/m3
156-60-5	trans-1,2-Dichloroethene	9.10	36.1		0.75	1.98	ug/m3
75-34-3	1,1-Dichloroethane	9.60	38.9		0.53	2.02	ug/m3
110-82-7	Cyclohexane	8.50	29.3		0.76	1.72	ug/m3
78-93-3	2-Butanone	8.30	24.5		0.35	1.47	ug/m3
56-23-5	Carbon Tetrachloride	9.50	59.8		0.060	0.19	ug/m3
156-59-2	cis-1,2-Dichloroethene	9.10	36.1		0.36	1.98	ug/m3
67-66-3	Chloroform	9.60	46.9		0.39	2.44	ug/m3
71-55-6	1,1,1-Trichloroethane	9.30	50.7		0.050	0.16	ug/m3
540-84-1	2,2,4-Trimethylpentane	8.10	37.8		0.47	2.34	ug/m3
71-43-2	Benzene	8.60	27.5		0.29	1.60	ug/m3
107-06-2	1,2-Dichloroethane	10.3	41.7		0.36	2.02	ug/m3
79-01-6	Trichloroethene	7.80	41.9		0.11	0.16	ug/m3
78-87-5	1,2-Dichloropropane	8.30	38.4		0.51	2.31	ug/m3
75-27-4	Bromodichloromethane	9.70	65.0		0.54	3.35	ug/m3
108-10-1	4-Methyl-2-Pentanone	7.20	29.5		0.37	2.05	ug/m3
108-88-3	Toluene	8.30	31.3		0.41	1.88	ug/m3
10061-02-6	t-1,3-Dichloropropene	8.50	38.6		0.27	2.27	ug/m3
10061-01-5	cis-1,3-Dichloropropene	8.30	37.7		0.27	2.27	ug/m3
79-00-5	1,1,2-Trichloroethane	9.00	49.1		0.38	2.73	ug/m3

Report of Analysis

Client:	GFE LLC	Date Collected:	
Project:	416 Clinton St Hempstead	Date Received:	
Client Sample ID:	VL0421ABS01	SDG No.:	Q1801
Lab Sample ID:	VL0421ABS01	Matrix:	Air
Analytical Method:	TO-15	Test:	TO-15
Sample Wt/Vol:	400	Units:	mL

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
VL042397.D	1		04/21/25 10:14	VL042125

CAS Number	Parameter	Conc. ppbv	Conc. ug/M3	Qualifier	MDL	LOQ / CRQL	Units
124-48-1	Dibromochloromethane	9.20	78.4		0.60	4.26	ug/m3
106-93-4	1,2-Dibromoethane	8.90	68.4		0.54	0.77	ug/m3
127-18-4	Tetrachloroethene	7.40	50.2		0.14	0.20	ug/m3
108-90-7	Chlorobenzene	8.90	41.0		0.37	2.30	ug/m3
100-41-4	Ethyl Benzene	9.00	39.1		0.52	2.17	ug/m3
179601-23-1	m/p-Xylene	18.2	79.0		0.91	4.34	ug/m3
95-47-6	o-Xylene	9.20	40.0		0.52	2.17	ug/m3
100-42-5	Styrene	8.60	36.6		0.47	2.13	ug/m3
75-25-2	Bromoform	8.80	91.0		0.62	5.17	ug/m3
79-34-5	1,1,2,2-Tetrachloroethane	8.40	57.7		0.070	0.21	ug/m3
95-49-8	2-Chlorotoluene	9.20	47.6		0.57	2.59	ug/m3
108-67-8	1,3,5-Trimethylbenzene	9.00	44.3		0.54	2.46	ug/m3
95-63-6	1,2,4-Trimethylbenzene	8.60	42.3		0.39	2.46	ug/m3
541-73-1	1,3-Dichlorobenzene	8.70	52.3		0.48	3.01	ug/m3
106-46-7	1,4-Dichlorobenzene	8.60	51.7		0.30	3.01	ug/m3
95-50-1	1,2-Dichlorobenzene	8.70	52.3		0.48	3.01	ug/m3
120-82-1	1,2,4-Trichlorobenzene	8.50	63.1		0.67	3.71	ug/m3
87-68-3	Hexachloro-1,3-Butadiene	8.50	90.7		0.96	5.33	ug/m3
106-99-0	1,3-Butadiene	9.00	19.9		0.29	1.11	ug/m3
91-20-3	Naphthalene	8.20	43.0		0.42	0.52	ug/m3
622-96-8	4-Ethyltoluene	9.00	44.3		0.59	2.46	ug/m3
110-54-3	Hexane	7.80	27.5		0.39	1.76	ug/m3
107-05-1	Allyl Chloride	9.10	28.5		0.47	1.57	ug/m3
123-91-1	1,4-Dioxane	7.40	26.7		0.76	1.80	ug/m3
80-62-6	Methyl Methacrylate	7.70	31.5		0.37	2.05	ug/m3
SURROGATES							
460-00-4	1-Bromo-4-Fluorobenzene	10.8			65 - 135	108%	SPK: 10
INTERNAL STANDARDS							
74-97-5	Bromochloromethane	118000			2.781		
540-36-3	1,4-Difluorobenzene	338000			3.949		
3114-55-4	Chlorobenzene-d5	294000			8.875		

Report of Analysis

Client:	GFE LLC	Date Collected:	
Project:	416 Clinton St Hempstead	Date Received:	
Client Sample ID:	VL0421ABS01	SDG No.:	Q1801
Lab Sample ID:	VL0421ABS01	Matrix:	Air
Analytical Method:	TO-15	Test:	TO-15
Sample Wt/Vol:	400	Units:	mL

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
VL042397.D	1		04/21/25 10:14	VL042125

CAS Number	Parameter	Conc. ppbv	Conc. ug/M3	Qualifier	MDL	LOQ / CRQL	Units
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U = Not Detected
 RL = Reporting Limit
 MDL = Method Detection Limit
 E = Value Exceeds Calibration Range
 D = Dilution

J = Estimated Value
 B = Analyte Found in Associated Method Blank
 N = Presumptive Evidence of a Compound
 * = Values outside of QC limits
 Q = indicates LCS control criteria did not meet requirements

Data Path : Z:\voasrv\HPCHEM1\MSVOA_L\Data\VL042125\
 Data File : VL042397.D
 Acq On : 21 Apr 2025 10:14
 Operator : SY/MD
 Sample : VL0421ABS01
 Misc : 400mL/MSVOA_L
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_L
 ClientSampleId :
 VL0421ABS01

Manual Integrations
 APPROVED

Reviewed By : Semsettin Yesilyurt 04/22/2025
 Supervised By : Mahesh Dadoda 04/22/2025

Quant Time: Apr 22 02:51:20 2025

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_L\methods\VL041725AIR.M

Quant Title : AIR ANALYSIS BY METHOD TO-15 Instrument: MSVOA_L Fri Aug 2025

QLast Update : Fri Apr 18 02:14:00 2025

Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) Bromochloromethane	2.781	49	118401	10.000	ppbv	0.00
33) 1,4-Difluorobenzene	3.949	114	337921	10.000	ppbv	0.00
55) Chlorobenzene-d5	8.875	117	293537	10.000	ppbv	0.00

System Monitoring Compounds

68) 1-Bromo-4-Fluorobenzene 10.371 95 260048 10.781 ppbv 0.00
 Spiked Amount 10.000 Range 65 - 135 Recovery = 107.800%

Target Compounds

						Qvalue
2) Dichlorodifluoromethane	1.496	85	212503	10.158	ppbv	98
3) Chlorodifluoromethane	1.473	51	175535	8.812	ppbv	95
4) Chloromethane	1.531	50	79791	9.638	ppbv	94
5) Vinyl Chloride	1.577	62	77905	8.645	ppbv	97
6) Bromomethane	1.664	94	35087	9.064	ppbv	87
7) Chloroethane	1.700	64	30814	9.455	ppbv	96
8) Dichlorotetrafluoroethane	1.551	85	182258	10.147	ppbv	94
9) Propene	1.480	41	69080	7.615	ppbv	97
10) Heptane	4.969	43	187072	7.340	ppbv	99
11) Trichlorofluoromethane	1.875	101	205641	10.337	ppbv	100
12) 1,1,2-Trichlorotrifluo...	2.133	101	156363	9.613	ppbv	98
13) Ethanol	1.716	45	4602	3.192	ppbv	92
14) Bromoethene	1.777	108	58938	9.613	ppbv	98
15) Acetone	1.832	43	168065	8.671	ppbv	98
16) 1,3-Butadiene	1.606	39	83577	9.015	ppbv	92
17) tert-Butyl alcohol	2.036	59	216440	9.011	ppbv	100
18) 1,1-Dichloroethene	2.030	96	71097	9.356	ppbv	95
19) Isopropyl Alcohol	1.881	45	96437	8.956	ppbv	93
20) Methylene Chloride	2.056	84	60417	7.959	ppbv	89
21) Allyl Chloride	2.091	41	122313	9.140	ppbv	97
22) trans-1,2-Dichloroethene	2.334	96	78976	9.128	ppbv	98
23) Vinyl Acetate	2.457	43	54610	6.961	ppbv #	96
24) 1,1-Dichloroethane	2.402	63	157093	9.559	ppbv	99
25) Ethyl Acetate	2.829	43	393203	8.228	ppbv	99
26) Hexane	2.820	57	151103	7.813	ppbv	99
27) Carbon Disulfide	2.146	76	194719	9.404	ppbv	98
28) Methyl tert-Butyl Ether	2.434	73	108007	8.524	ppbv	93
29) Chloroform	2.842	83	262639	9.608	ppbv	98
30) Cyclohexane	3.846	84	133153	8.517	ppbv	98
31) cis-1,2-Dichloroethene	2.713	61	166447	9.052	ppbv	95
32) 1,1,1-Trichloroethane	3.360	97	285156	9.266	ppbv	98
34) 2-Butanone	2.554	43	220966	8.268	ppbv	98
35) Carbon Tetrachloride	3.755	117	291941	9.546	ppbv	95
36) Benzene	3.648	78	312290	8.568	ppbv	97
37) 1,2-Dichloroethane	3.211	62	225346	10.301	ppbv #	96
38) Trichloroethene	4.542	130	123977	7.783	ppbv	98
39) 1,2-Dichloropropane	4.305	63	112058	8.250	ppbv #	90
40) 1,4-Dioxane	4.571	88	55511	7.437	ppbv #	93
41) Tetrahydrofuran	3.046	42	121497	7.797	ppbv	99
42) Bromodichloromethane	4.480	83	294127	9.661	ppbv	98
43) Methyl Methacrylate	4.872	69	125779	7.657	ppbv	93
44) 2,2,4-Trimethylpentane	4.629	57	508984	8.085	ppbv	95
45) t-1,3-Dichloropropene	6.399	75	142114	8.469	ppbv	93

Data Path : Z:\voasrv\HPCHEM1\MSVOA_L\Data\VL042125\
 Data File : VL042397.D
 Acq On : 21 Apr 2025 10:14
 Operator : SY/MD
 Sample : VL0421ABS01
 Misc : 400mL/MSVOA_L
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_L
 ClientSampleId :
 VL0421ABS01

Manual Integrations
 APPROVED

Reviewed By :Semsettin Yesilyurt 04/22/2025
 Supervised By :Mahesh Dadoda 04/22/2025

Quant Time: Apr 22 02:51:20 2025

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_L\methods\VL041725AIR.M

Quant Title : AIR ANALYSIS BY METHOD TO-15 Instrument: MSVOA_L Fri Aug 2

QLast Update : Fri Apr 18 02:14:00 2025

Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
46) cis-1,3-Dichloropropene	5.590	75	171445	8.348	ppbv #	84
47) 1,1,2-Trichloroethane	6.574	97	124092	8.967	ppbv	95
48) Dibromochloromethane	7.377	129	225553	9.226	ppbv	99
49) Bromoform	9.474	173	188764	8.830	ppbv	100
50) 4-Methyl-2-Pentanone	5.739	43	318982	7.182	ppbv	99
51) 2-Hexanone	7.429	43	266951	6.702	ppbv #	95
52) Tetrachloroethene	8.222	164	104899	7.446	ppbv	97
53) Toluene	6.920	91	361443	8.264	ppbv	99
54) 1,2-Dibromoethane	7.645	107	175799	8.877	ppbv	94
56) 1,1,1,2-Tetrachloroethane	8.917	131	169656	9.550	ppbv	97
57) Chlorobenzene	8.914	112	270733	8.868	ppbv	98
58) Ethyl Benzene	9.348	91	524070	9.011	ppbv	100
59) m/p-Xylene	9.545	91	852193m	18.189	ppbv	
60) o-Xylene	9.956	91	433062	9.179	ppbv	97
61) Styrene	9.866	104	194756	8.648	ppbv	98
62) Isopropylbenzene	10.533	105	641535	9.262	ppbv	97
63) 1,1,2,2-Tetrachloroethane	9.956	83	257428	8.420	ppbv	99
64) n-propylbenzene	10.982	120	155343	8.922	ppbv	96
65) tert-Butylbenzene	11.526	119	573929	9.015	ppbv	96
66) Benzyl Chloride	11.610	91	51025	7.294	ppbv	97
67) sec-Butylbenzene	11.762	105	802355	9.061	ppbv	99
69) p-Isopropyltoluene	11.921	119	685449	8.910	ppbv	97
70) n-Butylbenzene	12.277	91	723138	9.025	ppbv	98
71) 2-Chlorotoluene	10.895	91	500409	9.211	ppbv	97
72) 4-Ethyltoluene	11.122	105	519236	9.002	ppbv	98
73) 1,3,5-Trimethylbenzene	11.199	105	456404	8.963	ppbv	96
74) 1,2,4-Trimethylbenzene	11.533	105	509154	8.561	ppbv	91
75) 1,3-Dichlorobenzene	11.604	146	267483	8.652	ppbv	98
76) 1,4-Dichlorobenzene	11.669	146	267283	8.602	ppbv	98
77) 1,2-Dichlorobenzene	11.944	146	261316	8.693	ppbv	97
78) Hexachloro-1,3-Butadiene	13.879	225	259140	8.537	ppbv	99
79) Naphthalene	13.510	128	584987	8.225	ppbv	99
80) Naphthalene,2-methyl-	14.455	142	300148	7.405	ppbv	96
81) 1,2,4-Trichlorobenzene	13.436	180	258107	8.452	ppbv	99

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

Manual Integration Report

Sequence:	VL041725	Instrument	MSVOA_I
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Sample ID	File ID	Parameter	Review By	Review On	Supervised By	Supervised On	Reason
VSTDICCC010	VL042349.D	m/p-Xylene	SAM	4/18/2025 4:19:07 PM	MMDadoda	4/18/2025 10:45:20 PM	Peak Integrated by Software
VSTDICCC002	VL042350.D	1,4-Dioxane	SAM	4/18/2025 4:19:12 PM	MMDadoda	4/18/2025 10:45:17 PM	Peak Integrated by Software
VSTDICCC002	VL042350.D	cis-1,3-Dichloropropene	SAM	4/18/2025 4:19:12 PM	MMDadoda	4/18/2025 10:45:17 PM	Peak Integrated by Software
VSTDICCC002	VL042350.D	Ethanol	SAM	4/18/2025 4:19:12 PM	MMDadoda	4/18/2025 10:45:17 PM	Peak Integrated by Software
VSTDICCC002	VL042350.D	Heptane	SAM	4/18/2025 4:19:12 PM	MMDadoda	4/18/2025 10:45:17 PM	Peak Integrated by Software
VSTDICCC002	VL042350.D	m/p-Xylene	SAM	4/18/2025 4:19:12 PM	MMDadoda	4/18/2025 10:45:17 PM	Peak Integrated by Software
VSTDICCC002	VL042350.D	Trichloroethene	SAM	4/18/2025 4:19:12 PM	MMDadoda	4/18/2025 10:45:17 PM	Peak Integrated by Software
VSTDICCC001	VL042351.D	1,1,2-Trichloroethane	SAM	4/18/2025 4:19:17 PM	MMDadoda	4/18/2025 10:45:15 PM	Peak Integrated by Software
VSTDICCC001	VL042351.D	1,3-Butadiene	SAM	4/18/2025 4:19:17 PM	MMDadoda	4/18/2025 10:45:15 PM	Peak Integrated by Software
VSTDICCC001	VL042351.D	1,4-Dioxane	SAM	4/18/2025 4:19:17 PM	MMDadoda	4/18/2025 10:45:15 PM	Peak Integrated by Software
VSTDICCC001	VL042351.D	2,2,4-Trimethylpentane	SAM	4/18/2025 4:19:17 PM	MMDadoda	4/18/2025 10:45:15 PM	Peak Integrated by Software
VSTDICCC001	VL042351.D	cis-1,3-Dichloropropene	SAM	4/18/2025 4:19:17 PM	MMDadoda	4/18/2025 10:45:15 PM	Peak Integrated by Software
VSTDICCC001	VL042351.D	Cyclohexane	SAM	4/18/2025 4:19:17 PM	MMDadoda	4/18/2025 10:45:15 PM	Peak Integrated by Software

Manual Integration Report

Sequence:	VL041725	Instrument	MSVOA_I
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Sample ID	File ID	Parameter	Review By	Review On	Supervised By	Supervised On	Reason
VSTDICC001	VL042351.D	Ethanol	SAM	4/18/2025 4:19:17 PM	MMDadoda	4/18/2025 10:45:15 PM	Peak Integrated by Software
VSTDICC001	VL042351.D	Heptane	SAM	4/18/2025 4:19:17 PM	MMDadoda	4/18/2025 10:45:15 PM	Peak Integrated by Software
VSTDICC001	VL042351.D	m/p-Xylene	SAM	4/18/2025 4:19:17 PM	MMDadoda	4/18/2025 10:45:15 PM	Peak Integrated by Software
VSTDICC001	VL042351.D	Methyl Methacrylate	SAM	4/18/2025 4:19:17 PM	MMDadoda	4/18/2025 10:45:15 PM	Peak Integrated by Software
VSTDICC001	VL042351.D	t-1,3-Dichloropropene	SAM	4/18/2025 4:19:17 PM	MMDadoda	4/18/2025 10:45:15 PM	Peak Integrated by Software
VSTDICC0.5	VL042352.D	1,1,2-Trichloroethane	SAM	4/18/2025 4:20:33 PM	MMDadoda	4/18/2025 10:45:07 PM	Peak Integrated by Software
VSTDICC0.5	VL042352.D	1,2-Dichloropropane	SAM	4/18/2025 4:20:33 PM	MMDadoda	4/18/2025 10:45:07 PM	Peak Integrated by Software
VSTDICC0.5	VL042352.D	1,4-Dioxane	SAM	4/18/2025 4:20:33 PM	MMDadoda	4/18/2025 10:45:07 PM	Peak Integrated by Software
VSTDICC0.5	VL042352.D	4-Methyl-2-Pentanone	SAM	4/18/2025 4:20:33 PM	MMDadoda	4/18/2025 10:45:07 PM	Peak Integrated by Software
VSTDICC0.5	VL042352.D	Chlorobenzene	SAM	4/18/2025 4:20:33 PM	MMDadoda	4/18/2025 10:45:07 PM	Peak Integrated by Software
VSTDICC0.5	VL042352.D	cis-1,3-Dichloropropene	SAM	4/18/2025 4:20:33 PM	MMDadoda	4/18/2025 10:45:07 PM	Peak Integrated by Software
VSTDICC0.5	VL042352.D	Ethanol	SAM	4/18/2025 4:20:33 PM	MMDadoda	4/18/2025 10:45:07 PM	Peak Integrated by Software
VSTDICC0.5	VL042352.D	Heptane	SAM	4/18/2025 4:20:33 PM	MMDadoda	4/18/2025 10:45:07 PM	Peak Integrated by Software

Manual Integration Report

Sequence:	VL041725	Instrument	MSVOA_I
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Sample ID	File ID	Parameter	Review By	Review On	Supervised By	Supervised On	Reason
VSTDIC0.5	VL042352.D	m/p-Xylene	SAM	4/18/2025 4:20:33 PM	MMDadoda	4/18/2025 10:45:07 PM	Peak Integrated by Software
VSTDIC0.5	VL042352.D	Methyl Methacrylate	SAM	4/18/2025 4:20:33 PM	MMDadoda	4/18/2025 10:45:07 PM	Peak Integrated by Software
VSTDIC0.5	VL042352.D	t-1,3-Dichloropropene	SAM	4/18/2025 4:20:33 PM	MMDadoda	4/18/2025 10:45:07 PM	Peak Integrated by Software
VSTDIC0.5	VL042352.D	Tetrahydrofuran	SAM	4/18/2025 4:20:33 PM	MMDadoda	4/18/2025 10:45:07 PM	Peak Integrated by Software
VSTDIC0.1	VL042353.D	1,1,1-Trichloroethane	SAM	4/18/2025 4:20:37 PM	MMDadoda	4/18/2025 10:45:04 PM	Peak Integrated by Software
VSTDIC0.1	VL042353.D	1,1,2,2-Tetrachloroethane	SAM	4/18/2025 4:20:37 PM	MMDadoda	4/18/2025 10:45:04 PM	Peak Integrated by Software
VSTDIC0.1	VL042353.D	1,2-Dibromoethane	SAM	4/18/2025 4:20:37 PM	MMDadoda	4/18/2025 10:45:04 PM	Peak Integrated by Software
VSTDIC0.1	VL042353.D	Tetrachloroethene	SAM	4/18/2025 4:20:37 PM	MMDadoda	4/18/2025 10:45:04 PM	Peak Integrated by Software
VSTDIC0.1	VL042353.D	Trichloroethene	SAM	4/18/2025 4:20:37 PM	MMDadoda	4/18/2025 10:45:04 PM	Peak Integrated by Software
VSTDIC0.03	VL042354.D	1,1,1-Trichloroethane	SAM	4/18/2025 4:19:24 PM	MMDadoda	4/18/2025 10:45:02 PM	Peak Integrated by Software
VSTDIC0.03	VL042354.D	1,1,2,2-Tetrachloroethane	SAM	4/18/2025 4:19:24 PM	MMDadoda	4/18/2025 10:45:02 PM	Peak Integrated by Software
VSTDIC0.03	VL042354.D	Carbon Tetrachloride	SAM	4/18/2025 4:19:24 PM	MMDadoda	4/18/2025 10:45:02 PM	Peak Integrated by Software
VSTDIC0.03	VL042354.D	Tetrachloroethene	SAM	4/18/2025 4:19:24 PM	MMDadoda	4/18/2025 10:45:02 PM	Peak Integrated by Software

Manual Integration Report

Sequence:

vl041725

Instrument

MSVOA_I

Sample ID	File ID	Parameter	Review By	Review On	Supervised By	Supervised On	Reason
VSTDICC0.03	VL042354.D	Trichloroethene	SAM	4/18/2025 4:19:24 PM	MMDadoda	4/18/2025 10:45:02 PM	Peak Integrated by Software
VSTDICC015	VL042355.D	m/p-Xylene	SAM	4/18/2025 4:21:35 PM	MMDadoda	4/18/2025 10:44:59 PM	Peak Integrated by Software
VSTDICV010	VL042356.D	m/p-Xylene	SAM	4/18/2025 4:20:42 PM	MMDadoda	4/18/2025 10:44:57 PM	Peak Integrated by Software

Manual Integration Report

Sequence:	VL042125	Instrument	MSVOA_I
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Sample ID	File ID	Parameter	Review By	Review On	Supervised By	Supervised On	Reason
VSTDCCC010	VL042395.D	m/p-Xylene	SAM	4/22/2025 9:07:15 AM	MMDadoda	4/22/2025 1:59:41 PM	Peak Integrated by Software
VL0421ABS01	VL042397.D	m/p-Xylene	SAM	4/22/2025 9:07:22 AM	MMDadoda	4/22/2025 1:59:43 PM	Peak Integrated by Software
Q1801-01	VL042400.D	Ethanol	SAM	4/22/2025 9:07:29 AM	MMDadoda	4/22/2025 1:59:45 PM	Peak Integrated by Software
Q1801-01	VL042400.D	Heptane	SAM	4/22/2025 9:07:29 AM	MMDadoda	4/22/2025 1:59:45 PM	Peak Integrated by Software
Q1801-01	VL042400.D	m/p-Xylene	SAM	4/22/2025 9:07:29 AM	MMDadoda	4/22/2025 1:59:45 PM	Peak Integrated by Software
Q1801-02	VL042401.D	m/p-Xylene	SAM	4/22/2025 9:07:35 AM	MMDadoda	4/22/2025 1:59:47 PM	Peak Integrated by Software
Q1801-02	VL042401.D	Toluene	SAM	4/22/2025 9:07:35 AM	MMDadoda	4/22/2025 1:59:47 PM	Peak Integrated by Software
Q1844-02DUP	VL042403.D	Carbon Tetrachloride	SAM	4/22/2025 9:08:05 AM	MMDadoda	4/22/2025 1:59:51 PM	Peak Integrated by Software
Q1844-02DUP	VL042403.D	Cyclohexane	SAM	4/22/2025 9:08:05 AM	MMDadoda	4/22/2025 1:59:51 PM	Peak Integrated by Software
Q1844-02DUP	VL042403.D	Heptane	SAM	4/22/2025 9:08:05 AM	MMDadoda	4/22/2025 1:59:51 PM	Peak Integrated by Software
Q1844-02DUP	VL042403.D	m/p-Xylene	SAM	4/22/2025 9:08:05 AM	MMDadoda	4/22/2025 1:59:51 PM	Peak Integrated by Software
Q1844-02DUP	VL042403.D	Propene	SAM	4/22/2025 9:08:05 AM	MMDadoda	4/22/2025 1:59:51 PM	Peak Integrated by Software



284 Sheffield Street, Mountainside, New Jersey 07092, Phone : 908 789 8900, Fax : 908 789 8922

Manual Integration Report

Sequence:

VL042125

Instrument

MSVOA_I

Sample ID

File ID

Parameter

Review By

Review On

Supervised
By

Supervised On

Reason

Instrument ID: **MSVOA_L**

Daily Analysis Runlog For Sequence/QC Batch ID # VL041725

Review By	Semsettin Yesilyurt	Review On	4/18/2025 4:23:02 PM		
Supervise By	Maresh Dadoda	Supervise On	4/18/2025 10:46:12 PM		
SubDirectory	VL041725	HP Acquire Method	TO15AIR.M	HP Processing Method	VL041725AIR.M
STD. NAME		STD REF.#			
Tune/Reschk		AP2590			
Initial Calibration Stds		AP2598,AP2599,AP2600			
CCC		AP2598			
Internal Standard/PEM		AP2590			
ICV/I.BLK		AP2601			
Surrogate Standard					
MS/MSD Standard					
LCS Standard					

Sr#	SampleId	Data File Name	Date-Time	Operator	Status
1	BFB	VL042348.D	17 Apr 2025 07:50	SY/MD	Ok
2	VSTDICCC010	VL042349.D	17 Apr 2025 09:44	SY/MD	Ok,M
3	VSTDICCC002	VL042350.D	17 Apr 2025 10:17	SY/MD	Ok,M
4	VSTDICCC001	VL042351.D	17 Apr 2025 10:48	SY/MD	Ok,M
5	VSTDICCC0.5	VL042352.D	17 Apr 2025 11:19	SY/MD	Ok,M
6	VSTDICCC0.1	VL042353.D	17 Apr 2025 11:50	SY/MD	Ok,M
7	VSTDICCC0.03	VL042354.D	17 Apr 2025 12:22	SY/MD	Ok,M
8	VSTDICCC015	VL042355.D	17 Apr 2025 12:54	SY/MD	Ok,M
9	VSTDICV010	VL042356.D	17 Apr 2025 13:43	SY/MD	Ok,M
10	VL0417ABL01	VL042357.D	17 Apr 2025 14:23	SY/MD	Ok
11	VL0417ABS01	VL042358.D	17 Apr 2025 15:39	SY/MD	Ok,M
12	Q1665-15	VL042359.D	17 Apr 2025 16:12	SY/MD	Dilution
13	Q1665-15DL	VL042360.D	17 Apr 2025 16:44	SY/MD	Ok,M
14	Q1668-01	VL042361.D	17 Apr 2025 17:17	SY/MD	Dilution
15	Q1668-01DUP	VL042362.D	17 Apr 2025 17:51	SY/MD	Ok,M
16	Q1668-03	VL042363.D	17 Apr 2025 18:24	SY/MD	Dilution
17	Q1668-04	VL042364.D	17 Apr 2025 18:58	SY/MD	Ok,M
18	Q1668-02	VL042365.D	17 Apr 2025 19:31	SY/MD	Ok,M
19	Q1726-01	VL042366.D	17 Apr 2025 20:05	SY/MD	Dilution
20	Q1726-02	VL042367.D	17 Apr 2025 20:38	SY/MD	Dilution
21	Q1726-03	VL042368.D	17 Apr 2025 21:12	SY/MD	Dilution

Instrument ID: **MSVOA_L**

Daily Analysis Runlog For Sequence/QC Batch ID # VL041725

Review By	Semsettin Yesilyurt	Review On	4/18/2025 4:23:02 PM		
Supervise By	Maresh Dadoda	Supervise On	4/18/2025 10:46:12 PM		
SubDirectory	VL041725	HP Acquire Method	TO15AIR.M	HP Processing Method	VL041725AIR.M
STD. NAME		STD REF.#			
Tune/Reschk	AP2590				
Initial Calibration Stds	AP2598,AP2599,AP2600				
CCC	AP2598				
Internal Standard/PEM	AP2590				
ICV/I.BLK	AP2601				
Surrogate Standard					
MS/MSD Standard					
LCS Standard					

22	Q1726-04	VL042369.D	17 Apr 2025 21:45	SY/MD	Dilution
23	Q1726-05	VL042370.D	17 Apr 2025 22:19	SY/MD	Dilution
24	Q1726-06	VL042371.D	17 Apr 2025 22:52	SY/MD	Dilution
25	Q1726-07	VL042372.D	17 Apr 2025 23:26	SY/MD	Dilution
26	Q1726-07DL	VL042373.D	17 Apr 2025 23:57	SY/MD	Dilution
27	Q1668-01DL	VL042374.D	18 Apr 2025 00:28	SY/MD	Ok,M
28	Q1668-03DL	VL042375.D	18 Apr 2025 00:59	SY/MD	Ok,M
29	Q1726-01DL	VL042376.D	18 Apr 2025 01:30	SY/MD	Dilution
30	Q1726-02DL	VL042377.D	18 Apr 2025 02:01	SY/MD	Dilution
31	Q1726-03DL	VL042378.D	18 Apr 2025 02:32	SY/MD	Dilution
32	Q1726-04DL	VL042379.D	18 Apr 2025 03:03	SY/MD	Dilution
33	Q1726-05DL	VL042380.D	18 Apr 2025 03:34	SY/MD	Dilution
34	Q1726-06DL	VL042381.D	18 Apr 2025 04:05	SY/MD	Dilution

M : Manual Integration

Instrument ID: **MSVOA_L**

Daily Analysis Runlog For Sequence/QC Batch ID # VL042125

Review By	Semsettin Yesilyurt	Review On	4/22/2025 9:39:52 AM		
Supervise By	Maresh Dadoda	Supervise On	4/22/2025 2:00:33 PM		
SubDirectory	VL042125	HP Acquire Method	TO15AIR.M	HP Processing Method	VL041725AIR.M
STD. NAME		STD REF.#			
Tune/Reschk		AP2602			
Initial Calibration Stds		AP2598,AP2599,AP2600			
CCC		AP2598			
Internal Standard/PEM		AP2602			
ICV/I.BLK		AP2601			
Surrogate Standard					
MS/MSD Standard					
LCS Standard					

Sr#	SampleId	Data File Name	Date-Time	Operator	Status
1	BFB	VL042394.D	21 Apr 2025 07:58	SY/MD	Ok
2	VSTDCCC010	VL042395.D	21 Apr 2025 08:41	SY/MD	Ok,M
3	VL0421ABL01	VL042396.D	21 Apr 2025 09:28	SY/MD	Ok
4	VL0421ABS01	VL042397.D	21 Apr 2025 10:14	SY/MD	Ok,M
5	QC041625 CAN 10329	VL042398.D	21 Apr 2025 11:04	SY/MD	Not Ok
6	QC041625 CAN 10329	VL042399.D	21 Apr 2025 11:36	SY/MD	Ok
7	Q1801-01	VL042400.D	21 Apr 2025 12:07	SY/MD	Ok,M
8	Q1801-02	VL042401.D	21 Apr 2025 12:51	SY/MD	Ok,M
9	Q1844-02	VL042402.D	21 Apr 2025 13:38	SY/MD	Dilution
10	Q1844-02DUP	VL042403.D	21 Apr 2025 14:13	SY/MD	Ok,M
11	Q1844-02DL	VL042404.D	21 Apr 2025 14:44	SY/MD	Ok,M
12	Q1843-02	VL042405.D	21 Apr 2025 15:18	SY/MD	Ok,M
13	Q1843-02DL	VL042406.D	21 Apr 2025 15:49	SY/MD	Not Ok
14	Q1844-01	VL042407.D	21 Apr 2025 16:20	SY/MD	Dilution
15	Q1844-01DL	VL042408.D	21 Apr 2025 16:51	SY/MD	Dilution
16	Q1843-01	VL042409.D	21 Apr 2025 17:23	SY/MD	Ok,M
17	Q1843-01DL	VL042410.D	21 Apr 2025 17:54	SY/MD	Not Ok
18	Q1843-01DUP	VL042411.D	21 Apr 2025 19:16	SY/MD	Ok,M
19	QC041725 CA 27806	VL042412.D	21 Apr 2025 19:47	SY/MD	Not Ok
20	QC041725 CA 27806	VL042413.D	21 Apr 2025 20:18	SY/MD	Ok
21	Q1844-01DL2	VL042414.D	22 Apr 2025 07:47	SY/MD	Ok,M

M : Manual Integration

Instrument ID: MSVOA_L

Daily Analysis Runlog For Sequence/QC Batch ID # VL041725

Review By	Semsettin Yesilyurt	Review On	4/18/2025 4:23:02 PM
Supervise By	Maresh Dadoda	Supervise On	4/18/2025 10:46:12 PM
SubDirectory	VL041725	HP Acquire Method	TO15AIR.M HP Processing Method VL041725AIR.M

STD. NAME	STD REF.#
Tune/Reschk	AP2590
Initial Calibration Stds	AP2598,AP2599,AP2600
CCC	AP2598
Internal Standard/PEM	AP2590
ICV/I.BLK	AP2601
Surrogate Standard	
MS/MSD Standard	
LCS Standard	

Sr#	SampleID	ClientID	Data File Name	Date-Time	Comment	Operator	Status
1	BFB	BFB	VL042348.D	17 Apr 2025 07:50		SY/MD	Ok
2	VSTDICCC010	VSTDICCC010	VL042349.D	17 Apr 2025 09:44		SY/MD	Ok,M
3	VSTDICC002	VSTDICC002	VL042350.D	17 Apr 2025 10:17		SY/MD	Ok,M
4	VSTDICC001	VSTDICC001	VL042351.D	17 Apr 2025 10:48		SY/MD	Ok,M
5	VSTDICC0.5	VSTDICC0.5	VL042352.D	17 Apr 2025 11:19		SY/MD	Ok,M
6	VSTDICC0.1	VSTDICC0.1	VL042353.D	17 Apr 2025 11:50		SY/MD	Ok,M
7	VSTDICC0.03	VSTDICC0.03	VL042354.D	17 Apr 2025 12:22		SY/MD	Ok,M
8	VSTDICC015	VSTDICC015	VL042355.D	17 Apr 2025 12:54		SY/MD	Ok,M
9	VSTDICV010	ICVVL041725	VL042356.D	17 Apr 2025 13:43		SY/MD	Ok,M
10	VL0417ABL01	VL0417ABL01	VL042357.D	17 Apr 2025 14:23		SY/MD	Ok
11	VL0417ABS01	VL0417ABS01	VL042358.D	17 Apr 2025 15:39		SY/MD	Ok,M
12	Q1665-15	IA-B14-1-3242025	VL042359.D	17 Apr 2025 16:12	Need 10X	SY/MD	Dilution
13	Q1665-15DL	IA-B14-1-3242025DL	VL042360.D	17 Apr 2025 16:44		SY/MD	Ok,M
14	Q1668-01	IA-B1-4-3252025	VL042361.D	17 Apr 2025 17:17	Need 4X	SY/MD	Dilution
15	Q1668-01DUP	IA-B1-4-3252025DUP	VL042362.D	17 Apr 2025 17:51		SY/MD	Ok,M
16	Q1668-03	IA-B1-3-3252025	VL042363.D	17 Apr 2025 18:24	Need 4X	SY/MD	Dilution
17	Q1668-04	IA-B1-1-3252025	VL042364.D	17 Apr 2025 18:58		SY/MD	Ok,M
18	Q1668-02	IA-B1-2-3252025	VL042365.D	17 Apr 2025 19:31		SY/MD	Ok,M

Instrument ID: MSVOA_L

Daily Analysis Runlog For Sequence/QC Batch ID # VL041725

Review By	Semsettin Yesilyurt	Review On	4/18/2025 4:23:02 PM		
Supervise By	Mahesh Dadoda	Supervise On	4/18/2025 10:46:12 PM		
SubDirectory	VL041725	HP Acquire Method	TO15AIR.M	HP Processing Method	VL041725AIR.M
STD. NAME		STD REF.#			
Tune/Reschk		AP2590			
Initial Calibration Stds		AP2598,AP2599,AP2600			
CCC		AP2598			
Internal Standard/PEM		AP2590			
ICV/I.BLK		AP2601			
Surrogate Standard					
MS/MSD Standard					
LCS Standard					

19	Q1726-01	IA-DUP-3-03312025	VL042366.D	17 Apr 2025 20:05	Need 20X	SY/MD	Dilution
20	Q1726-02	IA-B2-6-03312025	VL042367.D	17 Apr 2025 20:38	Need 20X	SY/MD	Dilution
21	Q1726-03	IA-B2-2-03312025	VL042368.D	17 Apr 2025 21:12	Need 20X	SY/MD	Dilution
22	Q1726-04	IA-B2-5-03312025	VL042369.D	17 Apr 2025 21:45	Need 20X	SY/MD	Dilution
23	Q1726-05	IA-B2-1-03312025	VL042370.D	17 Apr 2025 22:19	Need 20X	SY/MD	Dilution
24	Q1726-06	IA-B2-3-03312025	VL042371.D	17 Apr 2025 22:52	Need 20X	SY/MD	Dilution
25	Q1726-07	IA-B2-4-03312025	VL042372.D	17 Apr 2025 23:26	Need 20X	SY/MD	Dilution
26	Q1726-07DL	IA-B2-4-03312025DL	VL042373.D	17 Apr 2025 23:57	Need 40X	SY/MD	Dilution
27	Q1668-01DL	IA-B1-4-3252025DL	VL042374.D	18 Apr 2025 00:28		SY/MD	Ok,M
28	Q1668-03DL	IA-B1-3-3252025DL	VL042375.D	18 Apr 2025 00:59		SY/MD	Ok,M
29	Q1726-01DL	IA-DUP-3-03312025DL	VL042376.D	18 Apr 2025 01:30	Need 40X	SY/MD	Dilution
30	Q1726-02DL	IA-B2-6-03312025DL	VL042377.D	18 Apr 2025 02:01	Need 40X	SY/MD	Dilution
31	Q1726-03DL	IA-B2-2-03312025DL	VL042378.D	18 Apr 2025 02:32	Need 100X	SY/MD	Dilution
32	Q1726-04DL	IA-B2-5-03312025DL	VL042379.D	18 Apr 2025 03:03	Need 40X	SY/MD	Dilution
33	Q1726-05DL	IA-B2-1-03312025DL	VL042380.D	18 Apr 2025 03:34	Need 100X	SY/MD	Dilution
34	Q1726-06DL	IA-B2-3-03312025DL	VL042381.D	18 Apr 2025 04:05	Need 100X	SY/MD	Dilution

M : Manual Integration

Instrument ID: MSVOA_L

Daily Analysis Runlog For Sequence/QC Batch ID # VL042125

Review By	Semsettin Yesilyurt	Review On	4/22/2025 9:39:52 AM
Supervise By	Maresh Dadoda	Supervise On	4/22/2025 2:00:33 PM
SubDirectory	VL042125	HP Acquire Method	TO15AIR.M
		HP Processing Method	VL041725AIR.M

STD. NAME	STD REF.#
Tune/Reschk	AP2602
Initial Calibration Stds	AP2598,AP2599,AP2600
CCC	AP2598
Internal Standard/PEM	AP2602
ICV/I.BLK	AP2601
Surrogate Standard	
MS/MSD Standard	
LCS Standard	

Sr#	SampleID	ClientID	Data File Name	Date-Time	Comment	Operator	Status
1	BFB	BFB	VL042394.D	21 Apr 2025 07:58		SY/MD	Ok
2	VSTDCCC010	VSTDCCC010	VL042395.D	21 Apr 2025 08:41		SY/MD	Ok,M
3	VL0421ABL01	VL0421ABL01	VL042396.D	21 Apr 2025 09:28		SY/MD	Ok
4	VL0421ABS01	VL0421ABS01	VL042397.D	21 Apr 2025 10:14		SY/MD	Ok,M
5	QC041625 CAN 10329	QC041625 CAN 10329	VL042398.D	21 Apr 2025 11:04	Contaminated	SY/MD	Not Ok
6	QC041625 CAN 10329	QC041625 CAN 10329	VL042399.D	21 Apr 2025 11:36		SY/MD	Ok
7	Q1801-01	SS2R	VL042400.D	21 Apr 2025 12:07		SY/MD	Ok,M
8	Q1801-02	SS3	VL042401.D	21 Apr 2025 12:51		SY/MD	Ok,M
9	Q1844-02	IA1	VL042402.D	21 Apr 2025 13:38	Need 5X	SY/MD	Dilution
10	Q1844-02DUP	IA1DUP	VL042403.D	21 Apr 2025 14:13		SY/MD	Ok,M
11	Q1844-02DL	IA1DL	VL042404.D	21 Apr 2025 14:44		SY/MD	Ok,M
12	Q1843-02	IA1	VL042405.D	21 Apr 2025 15:18		SY/MD	Ok,M
13	Q1843-02DL	IA1DL	VL042406.D	21 Apr 2025 15:49	Not required	SY/MD	Not Ok
14	Q1844-01	SV1	VL042407.D	21 Apr 2025 16:20	Need 40X	SY/MD	Dilution
15	Q1844-01DL	SV1DL	VL042408.D	21 Apr 2025 16:51	Need 600X	SY/MD	Dilution
16	Q1843-01	SV1	VL042409.D	21 Apr 2025 17:23		SY/MD	Ok,M
17	Q1843-01DL	SV1DL	VL042410.D	21 Apr 2025 17:54	Not required	SY/MD	Not Ok
18	Q1843-01DUP	SV1DUP	VL042411.D	21 Apr 2025 19:16		SY/MD	Ok,M

Instrument ID: MSVOA_L

Daily Analysis Runlog For Sequence/QC Batch ID # VL042125

Review By	Semsettin Yesilyurt	Review On	4/22/2025 9:39:52 AM
Supervise By	Maresh Dadoda	Supervise On	4/22/2025 2:00:33 PM
SubDirectory	VL042125	HP Acquire Method	TO15AIR.M
		HP Processing Method	VL041725AIR.M

STD. NAME	STD REF.#
Tune/Reschk	AP2602
Initial Calibration Stds	AP2598,AP2599,AP2600
CCC	AP2598
Internal Standard/PEM	AP2602
ICV/I.BLK	AP2601
Surrogate Standard	
MS/MSD Standard	
LCS Standard	

19	QC041725 CA 27806	QC041725 CA 27806	VL042412.D	21 Apr 2025 19:47	Contaminated	SY/MD	Not Ok
20	QC041725 CA 27806	QC041725 CA 27806	VL042413.D	21 Apr 2025 20:18		SY/MD	Ok
21	Q1844-01DL2	SV1DL2	VL042414.D	22 Apr 2025 07:47		SY/MD	Ok,M

M : Manual Integration

B2504002 - 3

Alliance Project No. :

New Jersey Department of Environmental Protection

Internal Chain of Custody

Instructions: Use 1 form for each 20 samples of aliquot

Laboratory: <u>Chemtech</u>		Location: <u>284 Sheffield Street, Mountinside, NJ 7092</u>	
NAME		Title: <u>Sample Custodian</u>	
Field Sample Seal No.: <u>Q1801</u>		Date Broken <u>4/14/2025</u>	
Case No.: <u>416 Clinton St Hempstead</u>		Analytical Parameter/Fraction <u>TO-15</u>	
		Military Time Seal Broken: <u>07:00:00</u>	

Sample No.	Aliquot/Extract No.	Sample No.	Aliquot/Extract No.
Q1801-01	SS2R		
Q1801-02	SS3		

Date	Time	Relinquished By	Received By	Purpose of Change of Custody
4/14	12:15	Signature:  Printed Name: <u>Douglas N.</u>	Signature:  Printed Name: <u>Samuel J. Jendryak</u>	
		Signature Printed Name	Signature Printed Name	
		Signature Printed Name	Signature Printed Name	
		Signature Printed Name	Signature Printed Name	
		Signature Printed Name	Signature Printed Name	
		Signature Printed Name	Signature Printed Name	
		Signature Printed Name	Signature Printed Name	
		Signature Printed Name	Signature Printed Name	
		Signature Printed Name	Signature Printed Name	
		Signature Printed Name	Signature Printed Name	

Distribution: White - Original (Sent With Report) Yellow - Contractor Archive Pink - Sample Custodian - Interim Copy

Analyst Signature:

[Signature]

Supervisor Signature:

2014

METHOD: TO-15

Pressure Gauge ID: A755971

Date	Sample Number	Canister #	Initial Pressure psia	Initial Pressure Hg	Final Pressure psia	Final Pressure Hg	Dilution Factor	Comment
04/08/15	Q1728-07	10467	9.8	-10.0				✓
	Q1728-08	10749	10.2	-9.2				
	Q1728-09	10722	9.6	-10.4				
	Q1728-10	10116	10.2	-9.2				
	Q1728-11	10466	9.6	-10.4				
✓	Q1728-12	10462	10.8	-8.0				✓
4/8/15	Q1757-01	10461	12.0	-5.5				✓
✓	Q1757-02	10258	11.6	-6.3				✓
04/09/15	Q1772-01	10439	12.2	-5.1				✓
"	Q1772-02	10285	12.4	-4.7				✓
04/09/15	Q1801-01	10311	13.6	-2.2				✓
	Q1801-02	10606	13.2	-3.1				
	Q1802-01	10329	12.4	-4.7				
	Q1802-02	10609	12.6	-4.3				
✓	Q1802-03	10334	13.2	-3.1				✓

Order ID : Q1801

Date	Bottle ID	Lab Sample Number	InComing Vacuum	Canister ID	Can Size	Reg ID	Reg Size	Batch Code	Certification ID
04/15/2025	B2504002	Q1801-02	-3.1	10606	6 L	10226		C2504001	VL042217.D Seq : VL040425 TUNE : VL042214.D ICAL : VL042177.D CCAL : VL042215.D MB : VL042216.D
04/15/2025	B2504002	Q1801-01	-2.2	10311	6 L	10648		C2504001	VL042217.D Seq : VL040425 TUNE : VL042214.D ICAL : VL042177.D CCAL : VL042215.D MB : VL042216.D

11301

22

CHEMTECH
284 Sheffield Street, Mountaintop, NJ 07092 P: (908) 789-8900 F: (908) 789-8922

Client Sample ID #: 552R
 Client Name: ATE
 Project Name: CLINTON
 Date: 4/12
 Analysis: TO-15
 Comments: _____
 CHEMTECH'S SAMPLE ID: _____
 Date Received: _____ Expiration: _____

Storage Location: L31
 Sample: Q1801-01
 Cust #: SS2R

002

90901

3-1

CHEMTECH
284 Sheffield Street, Mountaintop, NJ 07092 P: (908) 789-8900 F: (908) 789-8922

Client Sample ID #: 553
 Client Name: ATE
 Project Name: CLINTON ST
 Date: 4/12
 Analysis: TO-15
 Comments: _____
 CHEMTECH'S SAMPLE ID: _____
 Date Received: _____ Expiration: _____

Storage Location: L31
 Sample: Q1801-02
 Cust #: SS3

03504002