

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:	Q3757	OrderDate:	12/3/2025 9:25:00 AM
Client:	JACOBS Engineering Group, Inc.	Project:	Former Schlumberger STC PTC Site D3868221
Contact:	John Ynfante	Location:	--Select--,A11,VOA Lab

LabID	ClientID	Matrix	Test	Method	Sample Date	Prep Date	Anal Date	Received
Q3757-01	MW-14A-17.5-120325	Water	VOCMS Group3	8260-Low	12/03/25		12/05/25	12/03/25
Q3757-03	BR-03D-490-120325	Water	VOCMS Group3	8260-Low	12/03/25		12/05/25	12/03/25
Q3757-05	MW-14B-46-120325	Water	VOCMS Group3	8260-Low	12/03/25		12/05/25	12/03/25
Q3757-05RE	MW-14B-46-120325R E	Water	VOCMS Group3	8260-Low	12/03/25		12/08/25	12/03/25
Q3757-07	RMW-05B-90.5-12032 5	Water	VOCMS Group3	8260-Low	12/03/25		12/05/25	12/03/25
Q3757-09	OW-05B-72-120325	Water	VOCMS Group3	8260-Low	12/03/25		12/05/25	12/03/25
Q3757-09RE	OW-05B-72-120325R E	Water	VOCMS Group3	8260-Low	12/03/25		12/08/25	12/03/25
Q3757-11	EB01-120325	Water	VOCMS Group3	8260-Low	12/03/25		12/04/25	12/03/25
Q3757-12	TB01-120325	Water	VOCMS Group3	8260-Low	12/03/25		12/04/25	12/03/25

Hit Summary Sheet
SW-846

SDG No.: Q3757

Client: JACOBS Engineering Group, Inc.

Sample ID	Client ID	Matrix	Parameter	Concentration	C	MDL	RDL	Units
Client ID:	BR-03D-490-120325							
Q3757-03	BR-03D-490-12032	Water	Vinyl Chloride	3.90		0.26	1.00	ug/L
Q3757-03	BR-03D-490-12032	Water	1,1-Dichloroethene	48.1		0.23	1.00	ug/L
Q3757-03	BR-03D-490-12032	Water	1,1-Dichloroethane	6.40		0.23	1.00	ug/L
Q3757-03	BR-03D-490-12032	Water	cis-1,2-Dichloroethene	34.9		0.19	1.00	ug/L
Q3757-03	BR-03D-490-12032	Water	Benzene	0.23	J	0.15	1.00	ug/L
Q3757-03	BR-03D-490-12032	Water	1,2-Dichloroethane	3.70		0.22	1.00	ug/L
Q3757-03	BR-03D-490-12032	Water	Trichloroethene	35.1		0.090	1.00	ug/L
			Total Voc :	132				
			Total Concentration:	132				
Client ID:	RMW-05B-90.5-120325							
Q3757-07	RMW-05B-90.5-120	Water	1,1-Dichloroethane	1.50		0.23	1.00	ug/L
Q3757-07	RMW-05B-90.5-120	Water	cis-1,2-Dichloroethene	13.1		0.19	1.00	ug/L
Q3757-07	RMW-05B-90.5-120	Water	Trichloroethene	0.65	J	0.090	1.00	ug/L
			Total Voc :	15.3				
			Total Concentration:	15.3				
Client ID:	OW-05B-72-120325							
Q3757-09	OW-05B-72-12032	Water	1,1-Dichloroethene	0.85	J	0.23	1.00	ug/L
Q3757-09	OW-05B-72-12032	Water	cis-1,2-Dichloroethene	2.20		0.19	1.00	ug/L
Q3757-09	OW-05B-72-12032	Water	Trichloroethene	0.44	J	0.090	1.00	ug/L
			Total Voc :	3.49				
			Total Concentration:	3.49				
Client ID:	OW-05B-72-120325RE							
Q3757-09RE	OW-05B-72-12032	Water	1,1-Dichloroethene	0.60	J	0.23	1.00	ug/L
Q3757-09RE	OW-05B-72-12032	Water	cis-1,2-Dichloroethene	1.50		0.19	1.00	ug/L
Q3757-09RE	OW-05B-72-12032	Water	Trichloroethene	0.25	J	0.090	1.00	ug/L
			Total Voc :	2.35				
			Total Concentration:	2.35				



QC SUMMARY

Surrogate Summary

SDG No.: **Q3757**

Client: **JACOBS Engineering Group, Inc.**

Analytical Method: **SW8260-Low**

Lab Sample ID	Client ID	Parameter	Spike	Result	Recovery (%)	Qual	Limits (%)	
							Low	High
Q3757-01	MW-14A-17.5-120325	1,2-Dichloroethane-d4	50	52.4	105		70 (74)	130 (125)
		Dibromofluoromethane	50	49.5	99		70 (75)	130 (124)
		Toluene-d8	50	48.6	97		70 (86)	130 (113)
		4-Bromofluorobenzene	50	50.0	100		70 (77)	130 (121)
Q3757-03	BR-03D-490-120325	1,2-Dichloroethane-d4	50	55.1	110		70 (74)	130 (125)
		Dibromofluoromethane	50	51.8	104		70 (75)	130 (124)
		Toluene-d8	50	42.9	86		70 (86)	130 (113)
		4-Bromofluorobenzene	50	45.5	91		70 (77)	130 (121)
Q3757-05	MW-14B-46-120325	1,2-Dichloroethane-d4	50	47.2	94		70 (74)	130 (125)
		Dibromofluoromethane	50	40.9	82		70 (75)	130 (124)
		Toluene-d8	50	35.3	71		70 (86)	130 (113)
		4-Bromofluorobenzene	50	43.5	87		70 (77)	130 (121)
Q3757-05RE	MW-14B-46-120325RE	1,2-Dichloroethane-d4	50	57.7	115		70 (74)	130 (125)
		Dibromofluoromethane	50	38.4	77		70 (75)	130 (124)
		Toluene-d8	50	38.4	77		70 (86)	130 (113)
		4-Bromofluorobenzene	50	56.9	114		70 (77)	130 (121)
Q3757-07	RMW-05B-90.5-120325	1,2-Dichloroethane-d4	50	51.3	103		70 (74)	130 (125)
		Dibromofluoromethane	50	47.1	94		70 (75)	130 (124)
		Toluene-d8	50	43.9	88		70 (86)	130 (113)
		4-Bromofluorobenzene	50	45.8	92		70 (77)	130 (121)
Q3757-09	OW-05B-72-120325	1,2-Dichloroethane-d4	50	58.8	118		70 (74)	130 (125)
		Dibromofluoromethane	50	39.7	79		70 (75)	130 (124)
		Toluene-d8	50	38.2	76		70 (86)	130 (113)
		4-Bromofluorobenzene	50	45.0	90		70 (77)	130 (121)
Q3757-09RE	OW-05B-72-120325RE	1,2-Dichloroethane-d4	50	62.9	126		70 (74)	130 (125)
		Dibromofluoromethane	50	39.6	79		70 (75)	130 (124)
		Toluene-d8	50	47.0	94		70 (86)	130 (113)
		4-Bromofluorobenzene	50	56.1	112		70 (77)	130 (121)
Q3757-11	EB01-120325	1,2-Dichloroethane-d4	50	53.6	107		70 (74)	130 (125)
		Dibromofluoromethane	50	46.9	94		70 (75)	130 (124)
		Toluene-d8	50	44.4	89		70 (86)	130 (113)
		4-Bromofluorobenzene	50	49.5	99		70 (77)	130 (121)
Q3757-12	TB01-120325	1,2-Dichloroethane-d4	50	56.4	113		70 (74)	130 (125)
		Dibromofluoromethane	50	47.1	94		70 (75)	130 (124)
		Toluene-d8	50	45.0	90		70 (86)	130 (113)
		4-Bromofluorobenzene	50	52.4	105		70 (77)	130 (121)
VN1204WBL01	VN1204WBL01	1,2-Dichloroethane-d4	50	61.7	123		70 (74)	130 (125)
		Dibromofluoromethane	50	50.8	102		70 (75)	130 (124)
		Toluene-d8	50	46.7	93		70 (86)	130 (113)
		4-Bromofluorobenzene	50	49.9	100		70 (77)	130 (121)
VN1204WBS01	VN1204WBS01	1,2-Dichloroethane-d4	50	53.8	108		70 (74)	130 (125)
		Dibromofluoromethane	50	50.7	101		70 (75)	130 (124)
		Toluene-d8	50	51.7	103		70 (86)	130 (113)
		4-Bromofluorobenzene	50	54.4	109		70 (77)	130 (121)
VN1204WBSD0	VN1204WBSD01	1,2-Dichloroethane-d4	50	56.0	112		70 (74)	130 (125)
		Dibromofluoromethane	50	53.1	106		70 (75)	130 (124)
		Toluene-d8	50	52.7	105		70 (86)	130 (113)
		4-Bromofluorobenzene	50	56.3	113		70 (77)	130 (121)
VX1205WBL01	VX1205WBL01	1,2-Dichloroethane-d4	50	50.6	101		70 (74)	130 (125)
		Dibromofluoromethane	50	49.7	99		70 (75)	130 (124)

() = LABORATORY INHOUSE LIMIT

Surrogate Summary

SDG No.: Q3757

Client: JACOBS Engineering Group, Inc.

Analytical Method: SW8260-Low

Lab Sample ID	Client ID	Parameter	Spike	Result	Recovery (%)	Qual	Limits (%)	
							Low	High
VX1205WBL01	VX1205WBL01	Toluene-d8	50	50.8	102		70 (86)	130 (113)
		4-Bromofluorobenzene	50	60.5	121		70 (77)	130 (121)
VX1205WBS01	VX1205WBS01	1,2-Dichloroethane-d4	50	52.4	105		70 (74)	130 (125)
		Dibromofluoromethane	50	48.0	96		70 (75)	130 (124)
		Toluene-d8	50	49.6	99		70 (86)	130 (113)
		4-Bromofluorobenzene	50	52.5	105		70 (77)	130 (121)
VX1208WBL01	VX1208WBL01	1,2-Dichloroethane-d4	50	49.5	99		70 (74)	130 (125)
		Dibromofluoromethane	50	46.0	92		70 (75)	130 (124)
		Toluene-d8	50	45.2	90		70 (86)	130 (113)
		4-Bromofluorobenzene	50	55.8	112		70 (77)	130 (121)
VX1208WBS01	VX1208WBS01	1,2-Dichloroethane-d4	50	56.7	113		70 (74)	130 (125)
		Dibromofluoromethane	50	50.4	101		70 (75)	130 (124)
		Toluene-d8	50	50.6	101		70 (86)	130 (113)
		4-Bromofluorobenzene	50	58.8	118		70 (77)	130 (121)

Laboratory Control Sample/Laboratory Control Sample Duplicate Summary

SW-846

SDG No.:	<u>Q3757</u>	Analytical Method:	<u>SW8260-Low</u>
Client:	<u>JACOBS Engineering Group, Inc.</u>	Datafile :	<u>VN088443.D</u>

Lab Sample ID	Parameter	Spike	Result	Unit	Rec	RPD	Qual	Low	Limits High	RPD
VN1204WBS01	Vinyl chloride	20	20.3	ug/L	102			70 (65)	130 (117)	
	1,1-Dichloroethene	20	20.4	ug/L	102			70 (74)	130 (110)	
	1,1-Dichloroethane	20	19.6	ug/L	98			70 (78)	130 (112)	
	cis-1,2-Dichloroethene	20	19.0	ug/L	95			70 (77)	130 (110)	
	1,1,1-Trichloroethane	20	20.7	ug/L	104			70 (80)	130 (108)	
	Benzene	20	19.3	ug/L	97			70 (82)	130 (109)	
	1,2-Dichloroethane	20	20.8	ug/L	104			70 (80)	130 (115)	
	Trichloroethene	20	18.3	ug/L	92			70 (77)	130 (113)	
	1,1,2-Trichloroethane	20	19.9	ug/L	100			70 (83)	130 (112)	
	Tetrachloroethene	20	18.1	ug/L	91			70 (67)	130 (123)	

Laboratory Control Sample/Laboratory Control Sample Duplicate Summary

SW-846

SDG No.:	<u>Q3757</u>	Analytical Method:	<u>SW8260-Low</u>
Client:	<u>JACOBS Engineering Group, Inc.</u>	Datafile :	<u>VN088444.D</u>

Lab Sample ID	Parameter	Spike	Result	Unit	Rec	RPD	Qual	Low	Limits High	RPD
VN1204WBSD01	Vinyl chloride	20	19.0	ug/L	95	7		70 (65)	130 (117)	20 (19)
	1,1-Dichloroethene	20	19.5	ug/L	98	4		70 (74)	130 (110)	20 (20)
	1,1-Dichloroethane	20	19.7	ug/L	99	1		70 (78)	130 (112)	20 (20)
	cis-1,2-Dichloroethene	20	19.3	ug/L	97	2		70 (77)	130 (110)	20 (20)
	1,1,1-Trichloroethane	20	20.6	ug/L	103	1		70 (80)	130 (108)	20 (20)
	Benzene	20	19.3	ug/L	97	0		70 (82)	130 (109)	20 (15)
	1,2-Dichloroethane	20	21.6	ug/L	108	4		70 (80)	130 (115)	20 (20)
	Trichloroethene	20	18.7	ug/L	94	2		70 (77)	130 (113)	20 (15)
	1,1,2-Trichloroethane	20	20.7	ug/L	104	4		70 (83)	130 (112)	20 (20)
	Tetrachloroethene	20	17.3	ug/L	86	6		70 (67)	130 (123)	20 (15)

Laboratory Control Sample/Laboratory Control Sample Duplicate Summary

SW-846

SDG No.:	<u>Q3757</u>	Analytical Method:	<u>SW8260-Low</u>
Client:	<u>JACOBS Engineering Group, Inc.</u>	Datafile :	<u>VX048726.D</u>

Lab Sample ID	Parameter	Spike	Result	Unit	Rec	RPD	Qual	Low	Limits High	RPD
VX1205WBS01	Vinyl chloride	20	16.1	ug/L	81			70 (65)	130 (117)	
	1,1-Dichloroethene	20	19.4	ug/L	97			70 (74)	130 (110)	
	1,1-Dichloroethane	20	20.1	ug/L	101			70 (78)	130 (112)	
	cis-1,2-Dichloroethene	20	19.1	ug/L	96			70 (77)	130 (110)	
	1,1,1-Trichloroethane	20	20.0	ug/L	100			70 (80)	130 (108)	
	Benzene	20	18.7	ug/L	94			70 (82)	130 (109)	
	1,2-Dichloroethane	20	20.1	ug/L	101			70 (80)	130 (115)	
	Trichloroethene	20	19.0	ug/L	95			70 (77)	130 (113)	
	1,1,2-Trichloroethane	20	19.1	ug/L	96			70 (83)	130 (112)	
	Tetrachloroethene	20	17.6	ug/L	88			70 (67)	130 (123)	

Laboratory Control Sample/Laboratory Control Sample Duplicate Summary

SW-846

SDG No.:	<u>Q3757</u>	Analytical Method:	<u>SW8260-Low</u>
Client:	<u>JACOBS Engineering Group, Inc.</u>	Datafile :	<u>VX048745.D</u>

Lab Sample ID	Parameter	Spike	Result	Unit	Rec	RPD	Qual	Low	Limits High	RPD
VX1208WBS01	Vinyl chloride	20	19.8	ug/L	99			70 (65)	130 (117)	
	1,1-Dichloroethene	20	19.2	ug/L	96			70 (74)	130 (110)	
	1,1-Dichloroethane	20	20.8	ug/L	104			70 (78)	130 (112)	
	cis-1,2-Dichloroethene	20	19.9	ug/L	100			70 (77)	130 (110)	
	1,1,1-Trichloroethane	20	20.1	ug/L	101			70 (80)	130 (108)	
	Benzene	20	18.8	ug/L	94			70 (82)	130 (109)	
	1,2-Dichloroethane	20	21.4	ug/L	107			70 (80)	130 (115)	
	Trichloroethene	20	17.6	ug/L	88			70 (77)	130 (113)	
	1,1,2-Trichloroethane	20	21.7	ug/L	109			70 (83)	130 (112)	
	Tetrachloroethene	20	16.5	ug/L	83			70 (67)	130 (123)	

VOLATILE METHOD BLANK SUMMARY

Client ID

VN1204WBL01

Lab Name: Alliance

Contract: JAC005

Lab Code: ACE

SDG NO.: Q3757

Lab File ID: VN088439.D

Lab Sample ID: VN1204WBL01

Date Analyzed: 12/04/2025

Time Analyzed: 12:38

GC Column: RXI-624 ID: 0.25 (mm)

Heated Purge: (Y/N) N

Instrument ID: MSVOA_N

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

EPA SAMPLE NO.	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED
VN1204WBS01	VN1204WBS01	VN088443.D	12/04/2025
VN1204WBSD01	VN1204WBSD01	VN088444.D	12/04/2025
TB01-120325	Q3757-12	VN088449.D	12/04/2025
EB01-120325	Q3757-11	VN088450.D	12/04/2025

COMMENTS: _____

VOLATILE METHOD BLANK SUMMARY

Client ID

VX1205WBL01

Lab Name: Alliance

Contract: JAC005

Lab Code: ACE

SDG NO.: Q3757

Lab File ID: VX048725.D

Lab Sample ID: VX1205WBL01

Date Analyzed: 12/05/2025

Time Analyzed: 10:34

GC Column: DB-624UI ID: 0.18 (mm)

Heated Purge: (Y/N) N

Instrument ID: MSVOA_X

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

EPA SAMPLE NO.	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED
VX1205WBS01	VX1205WBS01	VX048726.D	12/05/2025
MW-14A-17.5-120325	Q3757-01	VX048731.D	12/05/2025
BR-03D-490-120325	Q3757-03	VX048732.D	12/05/2025
MW-14B-46-120325	Q3757-05	VX048733.D	12/05/2025
RMW-05B-90.5-120325	Q3757-07	VX048734.D	12/05/2025
OW-05B-72-120325	Q3757-09	VX048735.D	12/05/2025

COMMENTS: _____

VOLATILE METHOD BLANK SUMMARY

Client ID

VX1208WBL01

Lab Name: Alliance

Contract: JAC005

Lab Code: ACE

SDG NO.: Q3757

Lab File ID: VX048744.D

Lab Sample ID: VX1208WBL01

Date Analyzed: 12/08/2025

Time Analyzed: 11:15

GC Column: DB-624UI ID: 0.18 (mm)

Heated Purge: (Y/N) N

Instrument ID: MSVOA_X

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

EPA SAMPLE NO.	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED
VX1208WBS01	VX1208WBS01	VX048745.D	12/08/2025
MW-14B-46-120325RE	Q3757-05RE	VX048747.D	12/08/2025
OW-05B-72-120325RE	Q3757-09RE	VX048748.D	12/08/2025

COMMENTS: _____



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

VOLATILE ORGANIC INSTRUMENT PERFORMANCE CHECK
BROMOFLUOROBENZENE (BFB)

Lab Name: Alliance

Contract: JAC005

Lab Code: ACE

SDG NO.: Q3757

Lab File ID: VN088343.D

BFB Injection Date: 11/24/2025

Instrument ID: MSVOA_N

BFB Injection Time: 08:17

GC Column: RXI-624 ID: 0.25 (mm)

Heated Purge: Y/N N

m/e	ION ABUNDANCE CRITERIA	% RELATIVE ABUNDANCE
50	15.0 - 40.0% of mass 95	23.6
75	30.0 - 60.0% of mass 95	56.6
95	Base Peak, 100% relative abundance	100
96	5.0 - 9.0% of mass 95	6.9
173	Less than 2.0% of mass 174	1 (1.4) 1
174	50.0 - 100.0% of mass 95	70.8
175	5.0 - 9.0% of mass 174	6 (8.4) 1
176	95.0 - 101.0% of mass 174	68.3 (96.4) 1
177	5.0 - 9.0% of mass 176	4.3 (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:

CLIENT ID	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED	TIME ANALYZED
VSTDIC001	VSTDIC001	VN088344.D	11/24/2025	12:38
VSTDIC005	VSTDIC005	VN088345.D	11/24/2025	13:11
VSTDIC020	VSTDIC020	VN088346.D	11/24/2025	13:32
VSTDIC050	VSTDIC050	VN088347.D	11/24/2025	13:53
VSTDIC100	VSTDIC100	VN088348.D	11/24/2025	14:14
VSTDIC150	VSTDIC150	VN088349.D	11/24/2025	14:35



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

VOLATILE ORGANIC INSTRUMENT PERFORMANCE CHECK
BROMOFLUOROBENZENE (BFB)

Lab Name: Alliance

Contract: JACO05

Lab Code: ACE

SDG NO.: Q3757

Lab File ID: VN088436.D

BFB Injection Date: 12/04/2025

Instrument ID: MSVOA_N

BFB Injection Time: 11:03

GC Column: RXI-624 ID: 0.25 (mm)

Heated Purge: Y/N N

m/e	ION ABUNDANCE CRITERIA	% RELATIVE ABUNDANCE
50	15.0 - 40.0% of mass 95	23.6
75	30.0 - 60.0% of mass 95	59.8
95	Base Peak, 100% relative abundance	100
96	5.0 - 9.0% of mass 95	7
173	Less than 2.0% of mass 174	0.5 (0.7) 1
174	50.0 - 100.0% of mass 95	70.2
175	5.0 - 9.0% of mass 174	5.8 (8.3) 1
176	95.0 - 101.0% of mass 174	67.6 (96.3) 1
177	5.0 - 9.0% of mass 176	4.3 (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:

CLIENT ID	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED	TIME ANALYZED
VSTDCCC050	VSTDCCC050	VN088437.D	12/04/2025	11:44
VN1204WBL01	VN1204WBL01	VN088439.D	12/04/2025	12:38
VN1204WBS01	VN1204WBS01	VN088443.D	12/04/2025	14:12
VN1204WBSD01	VN1204WBSD01	VN088444.D	12/04/2025	14:45
TB01-120325	Q3757-12	VN088449.D	12/04/2025	16:28
EB01-120325	Q3757-11	VN088450.D	12/04/2025	16:48



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

VOLATILE ORGANIC INSTRUMENT PERFORMANCE CHECK
BROMOFLUOROBENZENE (BFB)

Lab Name: Alliance

Contract: JACO05

Lab Code: ACE

SDG NO.: Q3757

Lab File ID: VX048713.D

BFB Injection Date: 12/04/2025

Instrument ID: MSVOA_X

BFB Injection Time: 08:15

GC Column: DB-624UI ID: 0.18 (mm)

Heated Purge: Y/N N

m/e	ION ABUNDANCE CRITERIA	% RELATIVE ABUNDANCE
50	15.0 - 40.0% of mass 95	19.6
75	30.0 - 60.0% of mass 95	51.6
95	Base Peak, 100% relative abundance	100
96	5.0 - 9.0% of mass 95	6.4
173	Less than 2.0% of mass 174	1 (1.3) 1
174	50.0 - 100.0% of mass 95	76.2
175	5.0 - 9.0% of mass 174	5.9 (7.8) 1
176	95.0 - 101.0% of mass 174	74.4 (97.7) 1
177	5.0 - 9.0% of mass 176	4.9 (6.6) 2

1-Value is % mass 174

2-Value is % mass 176

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

CLIENT ID	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED	TIME ANALYZED
VSTDIC001	VSTDIC001	VX048714.D	12/04/2025	08:47
VSTDIC005	VSTDIC005	VX048715.D	12/04/2025	09:16
VSTDIC020	VSTDIC020	VX048716.D	12/04/2025	09:36
VSTDIC050	VSTDIC050	VX048717.D	12/04/2025	09:57
VSTDIC100	VSTDIC100	VX048718.D	12/04/2025	10:17
VSTDIC150	VSTDIC150	VX048719.D	12/04/2025	10:38



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

VOLATILE ORGANIC INSTRUMENT PERFORMANCE CHECK
BROMOFLUOROBENZENE (BFB)

Lab Name: Alliance
Lab Code: ACE
Lab File ID: VX048722.D
Instrument ID: MSVOA_X
GC Column: DB-624UI ID: 0.18 (mm)

Contract: JACO05
SDG NO.: Q3757
BFB Injection Date: 12/05/2025
BFB Injection Time: 08:14
Heated Purge: Y/N N

m/e	ION ABUNDANCE CRITERIA	% RELATIVE ABUNDANCE
50	15.0 - 40.0% of mass 95	16.3
75	30.0 - 60.0% of mass 95	48.5
95	Base Peak, 100% relative abundance	100
96	5.0 - 9.0% of mass 95	6.6
173	Less than 2.0% of mass 174	0.9 (1.2) 1
174	50.0 - 100.0% of mass 95	80.6
175	5.0 - 9.0% of mass 174	5.9 (7.3) 1
176	95.0 - 101.0% of mass 174	77.7 (96.4) 1
177	5.0 - 9.0% of mass 176	5.4 (7) 2

1-Value is % mass 174

2-Value is % mass 176

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

CLIENT ID	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED	TIME ANALYZED
VSTDCCC050	VSTDCCC050	VX048723.D	12/05/2025	08:45
VX1205WBL01	VX1205WBL01	VX048725.D	12/05/2025	10:34
VX1205WBS01	VX1205WBS01	VX048726.D	12/05/2025	11:31
MW-14A-17.5-120325	Q3757-01	VX048731.D	12/05/2025	13:32
BR-03D-490-120325	Q3757-03	VX048732.D	12/05/2025	13:53
MW-14B-46-120325	Q3757-05	VX048733.D	12/05/2025	14:13
RMW-05B-90.5-120325	Q3757-07	VX048734.D	12/05/2025	14:33
OW-05B-72-120325	Q3757-09	VX048735.D	12/05/2025	14:54



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

VOLATILE ORGANIC INSTRUMENT PERFORMANCE CHECK
BROMOFLUOROBENZENE (BFB)

Lab Name: Alliance

Contract: JACO05

Lab Code: ACE

SDG NO.: Q3757

Lab File ID: VX048741.D

BFB Injection Date: 12/08/2025

Instrument ID: MSVOA_X

BFB Injection Time: 08:17

GC Column: DB-624UI ID: 0.18 (mm)

Heated Purge: Y/N N

m/e	ION ABUNDANCE CRITERIA	% RELATIVE ABUNDANCE
50	15.0 - 40.0% of mass 95	19.5
75	30.0 - 60.0% of mass 95	51.5
95	Base Peak, 100% relative abundance	100
96	5.0 - 9.0% of mass 95	6.2
173	Less than 2.0% of mass 174	1 (1.2) 1
174	50.0 - 100.0% of mass 95	80.1
175	5.0 - 9.0% of mass 174	6 (7.5) 1
176	95.0 - 101.0% of mass 174	76.4 (95.3) 1
177	5.0 - 9.0% of mass 176	5.3 (7) 2

1-Value is % mass 174

2-Value is % mass 176

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

CLIENT ID	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED	TIME ANALYZED
VSTDCCC050	VSTDCCC050	VX048742.D	12/08/2025	09:20
VX1208WBL01	VX1208WBL01	VX048744.D	12/08/2025	11:15
VX1208WBS01	VX1208WBS01	VX048745.D	12/08/2025	11:44
MW-14B-46-120325RE	Q3757-05RE	VX048747.D	12/08/2025	12:33
OW-05B-72-120325RE	Q3757-09RE	VX048748.D	12/08/2025	12:53

VOLATILE INTERNAL STANDARD AREA AND RT SUMMARY

Lab Name: Alliance Contract: JACO05
 Lab Code: ACE SDG NO.: Q3757
 Lab File ID: VN088437.D Date Analyzed: 12/04/2025
 Instrument ID: MSVOA_N Time Analyzed: 11:44
 GC Column: RXI-624 ID: 0.25 (mm) Heated Purge: (Y/N) N

	IS1 AREA #	RT #	IS2 AREA #	RT #	IS3 AREA #	RT #
12 HOUR STD	259230	8.21	454552	9.08	431157	11.84
UPPER LIMIT	518460	8.706	909104	9.577	862314	12.341
LOWER LIMIT	129615	7.706	227276	8.577	215579	11.341
EPA SAMPLE NO.						
EB01-120325	201753	8.21	451562	9.08	439722	11.84
TB01-120325	198228	8.21	454339	9.08	441078	11.84
VN1204WBL01	236600	8.21	556948	9.08	555675	11.84
VN1204WBS01	303745	8.21	559564	9.08	491842	11.84
VN1204WBSD01	290727	8.21	529277	9.08	469866	11.84

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

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

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

VOLATILE INTERNAL STANDARD AREA AND RT SUMMARY

Lab Name: Alliance Contract: JAC005
 Lab Code: ACE SDG NO.: Q3757
 Lab File ID: VN088437.D Date Analyzed: 12/04/2025
 Instrument ID: MSVOA_N Time Analyzed: 11:44
 GC Column: RXI-624 ID: 0.25 (mm) Heated Purge: (Y/N) N

	IS4 AREA #	RT #				
12 HOUR STD	210277	13.765				
UPPER LIMIT	420554	14.265				
LOWER LIMIT	105139	13.265				
EPA SAMPLE NO.						
EB01-120325	210586	13.77				
TB01-120325	211349	13.77				
VN1204WBL01	258091	13.77				
VN1204WBS01	237731	13.77				
VN1204WBSD01	229597	13.77				

IS4 = 1,4-Dichlorobenzene-d4

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

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

VOLATILE INTERNAL STANDARD AREA AND RT SUMMARY

Lab Name: Alliance Contract: JAC005
 Lab Code: ACE SDG NO.: Q3757
 Lab File ID: VX048723.D Date Analyzed: 12/05/2025
 Instrument ID: MSVOA_X Time Analyzed: 08:45
 GC Column: DB-624UI ID: 0.18 (mm) Heated Purge: (Y/N) N

	IS1 AREA #	RT #	IS2 AREA #	RT #	IS3 AREA #	RT #
12 HOUR STD	218600	5.53	333007	6.74	285113	10.04
UPPER LIMIT	437200	6.025	666014	7.239	570226	10.537
LOWER LIMIT	109300	5.025	166504	6.239	142557	9.537
EPA SAMPLE NO.						
MW-14A-17.5-120325	176619	5.53	302604	6.74	274042	10.04
BR-03D-490-120325	163180	5.54	280929	6.75	230844	10.04
MW-14B-46-120325	161591	5.54	286549	6.75	220300	10.04
RMW-05B-90.5-120325	132392	5.54	222932	6.75	207009	10.04
OW-05B-72-120325	129477	5.54	275147	6.75	208806	10.04
VX1205WBL01	193187	5.53	291151	6.74	294164	10.04
VX1205WBS01	179194	5.53	293931	6.74	269397	10.04

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

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

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

VOLATILE INTERNAL STANDARD AREA AND RT SUMMARY

Lab Name: Alliance Contract: JAC005
 Lab Code: ACE SDG NO.: Q3757
 Lab File ID: VX048723.D Date Analyzed: 12/05/2025
 Instrument ID: MSVOA_X Time Analyzed: 08:45
 GC Column: DB-624UI ID: 0.18 (mm) Heated Purge: (Y/N) N

	IS4 AREA #	RT #				
12 HOUR STD	141895	12				
UPPER LIMIT	283790	12.5				
LOWER LIMIT	70947.5	11.5				
EPA SAMPLE NO.						
MW-14A-17.5-120325	141373	12.01				
BR-03D-490-120325	110489	12.01				
MW-14B-46-120325	131055	12.01				
RMW-05B-90.5-120325	102583	12.01				
OW-05B-72-120325	103213	12.01				
VX1205WBL01	130501	12.01				
VX1205WBS01	132992	12.01				

IS4 = 1,4-Dichlorobenzene-d4

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

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

VOLATILE INTERNAL STANDARD AREA AND RT SUMMARY

Lab Name: Alliance Contract: JACO05
 Lab Code: ACE SDG NO.: Q3757
 Lab File ID: VX048742.D Date Analyzed: 12/08/2025
 Instrument ID: MSVOA_X Time Analyzed: 09:20
 GC Column: DB-624UI ID: 0.18 (mm) Heated Purge: (Y/N) N

	IS1 AREA #	RT #	IS2 AREA #	RT #	IS3 AREA #	RT #
12 HOUR STD	174813	5.53	271670	6.74	259893	10.04
UPPER LIMIT	349626	6.031	543340	7.239	519786	10.537
LOWER LIMIT	87406.5	5.031	135835	6.239	129947	9.537
EPA SAMPLE NO.						
MW-14B-46-120325RE	124903	5.54	299763	6.75	268328	10.04
OW-05B-72-120325RE	169111	5.54	409831	6.74	413547	10.04
VX1208WBL01	173343	5.53	340953	6.74	370451	10.04
VX1208WBS01	178446	5.53	298973	6.74	311271	10.04

IS1 = Pentafluorobenzene

IS2 = 1,4-Difluorobenzene

IS3 = Chlorobenzene-d5

AREA UPPER LIMIT = +100% of internal standard area

AREA LOWER LIMIT = -50% of internal standard area

RT UPPER LIMIT = +0.50 minutes of internal standard RT

RT LOWER LIMIT = -0.50 minutes of internal standard RT

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

* Values outside of QC limits.

VOLATILE INTERNAL STANDARD AREA AND RT SUMMARY

Lab Name: Alliance Contract: JAC005
 Lab Code: ACE SDG NO.: Q3757
 Lab File ID: VX048742.D Date Analyzed: 12/08/2025
 Instrument ID: MSVOA_X Time Analyzed: 09:20
 GC Column: DB-624UI ID: 0.18 (mm) Heated Purge: (Y/N) N

	IS4 AREA #	RT #				
12 HOUR STD	132353	12				
UPPER LIMIT	264706	12.5				
LOWER LIMIT	66176.5	11.5				
EPA SAMPLE NO.						
MW-14B-46-120325RE	139875	12.01				
OW-05B-72-120325RE	197728	12.01				
VX1208WBL01	179749	12.01				
VX1208WBS01	157846	12.00				

IS4 = 1,4-Dichlorobenzene-d4

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

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



SAMPLE DATA



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

Report of Analysis

Client: JACOBS Engineering Group, Inc.
Project: Former Schlumberger STC PTC Site D3868221
Client Sample ID: MW-14A-17.5-120325
Lab Sample ID: Q3757-01
Analytical Method: 8260D
Sample Wt/Vol: 5 mL

Level: LOW
Final Vol: 5000 uL

Date Collected: 12/03/25
Date Received: 12/03/25
SDG No.: Q3757
Matrix: Water
% Solid: 0
Test: VOCMS Group3

CAS Number	Parameter	Conc.	Qua.	DF	MDL	LOQ / CRQL	Units	Date Ana.	BatchID
TARGETS									
75-01-4	Vinyl Chloride	0.26	U	1	0.26	1.00	ug/L	12/05/25 13:32	VX120525
75-35-4	1,1-Dichloroethene	0.23	U	1	0.23	1.00	ug/L	12/05/25 13:32	VX120525
75-34-3	1,1-Dichloroethane	0.23	U	1	0.23	1.00	ug/L	12/05/25 13:32	VX120525
156-59-2	cis-1,2-Dichloroethene	0.19	U	1	0.19	1.00	ug/L	12/05/25 13:32	VX120525
71-55-6	1,1,1-Trichloroethane	0.20	U	1	0.20	1.00	ug/L	12/05/25 13:32	VX120525
71-43-2	Benzene	0.15	U	1	0.15	1.00	ug/L	12/05/25 13:32	VX120525
107-06-2	1,2-Dichloroethane	0.22	U	1	0.22	1.00	ug/L	12/05/25 13:32	VX120525
79-01-6	Trichloroethene	0.090	U	1	0.090	1.00	ug/L	12/05/25 13:32	VX120525
79-00-5	1,1,2-Trichloroethane	0.21	U	1	0.21	1.00	ug/L	12/05/25 13:32	VX120525
127-18-4	Tetrachloroethene	0.23	U	1	0.23	1.00	ug/L	12/05/25 13:32	VX120525
SURROGATES									
17060-07-0	1,2-Dichloroethane-d4	52.4			70 (74) - 130 (125)	105%	SPK: 50		
1868-53-7	Dibromofluoromethane	49.5			70 (75) - 130 (124)	99%	SPK: 50		
2037-26-5	Toluene-d8	48.7			70 (86) - 130 (113)	97%	SPK: 50		
460-00-4	4-Bromofluorobenzene	50.0			70 (77) - 130 (121)	100%	SPK: 50		
INTERNAL STANDARDS									
		Area Count							
363-72-4	Pentafluorobenzene	177000							
540-36-3	1,4-Difluorobenzene	303000							
3114-55-4	Chlorobenzene-d5	274000							
3855-82-1	1,4-Dichlorobenzene-d4	141000							

U = Not Detected

LOQ = Limit of Quantitation

MDL = Method Detection Limit

LOD = Limit of Detection

E = Value Exceeds Calibration Range

Q = indicates LCS control criteria did not meet requirements

M = MS/MSD acceptance criteria did not meet requirements

J = Estimated Value

B = Analyte Found in Associated Method Blank

N = Presumptive Evidence of a Compound

* = Values outside of QC limits

D = Dilution

() = Laboratory InHouse Limit

A = Aldol-Condensation Reaction Products

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120525\
 Data File : VX048731.D
 Acq On : 05 Dec 2025 13:32
 Operator : JC/MD
 Sample : Q3757-01
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-14A-17.5-120325

Quant Time: Dec 05 22:03:35 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X120425W.M
 Quant Title : SW846 8260
 QLast Update : Fri Dec 05 03:27:34 2025
 Response via : Initial Calibration

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

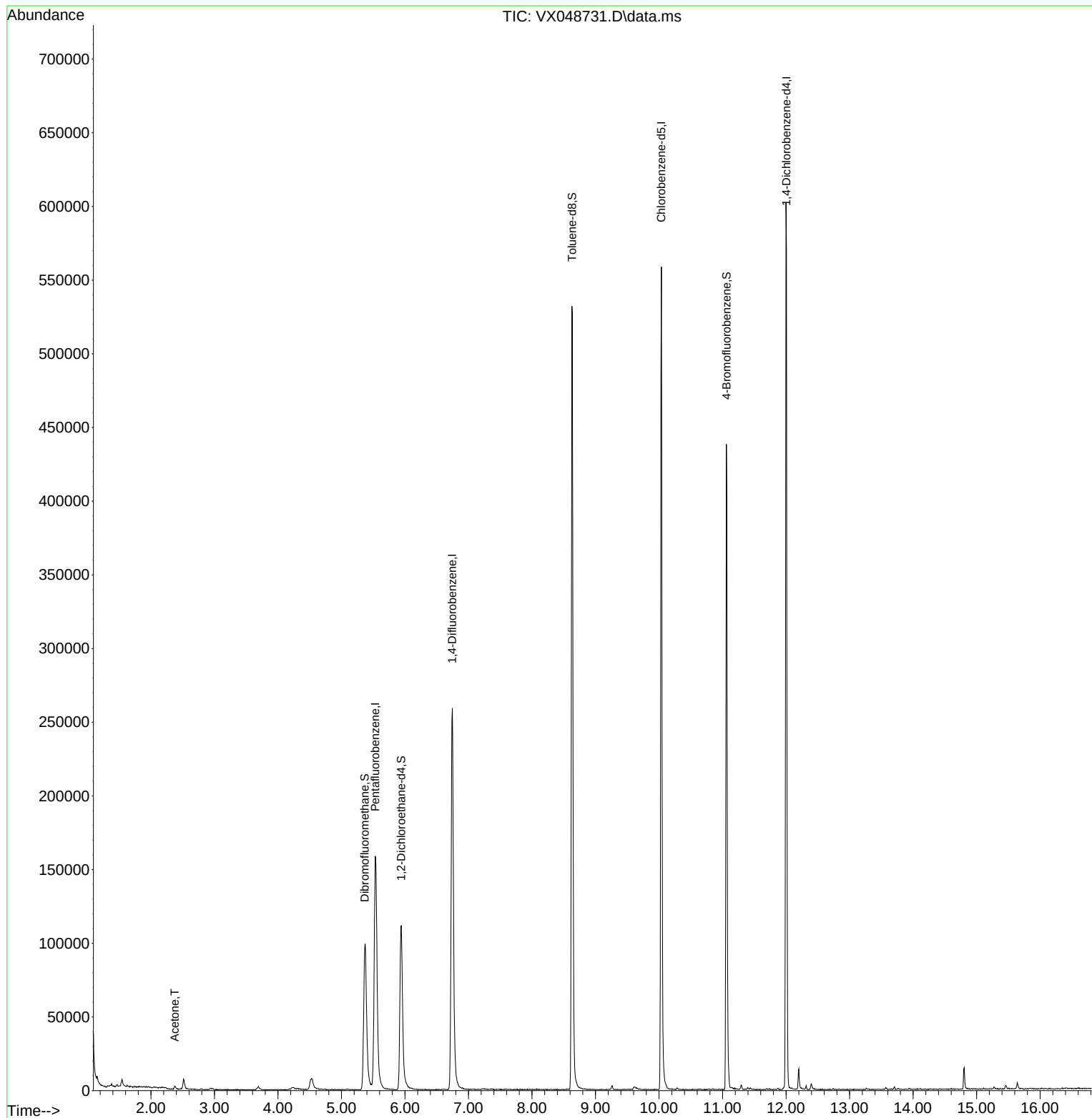
Internal Standards						
1) Pentafluorobenzene	5.531	168	176619	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.744	114	302604	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.037	117	274042	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.006	152	141373	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.940	65	124177	52.383	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	104.760%
35) Dibromofluoromethane	5.367	113	103142	49.503	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	99.000%
50) Toluene-d8	8.628	98	355343	48.654	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	97.300%
62) 4-Bromofluorobenzene	11.061	95	125871	49.979	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	99.960%
Target Compounds						
16) Acetone	2.373	43	3056	3.368	ug/l	95

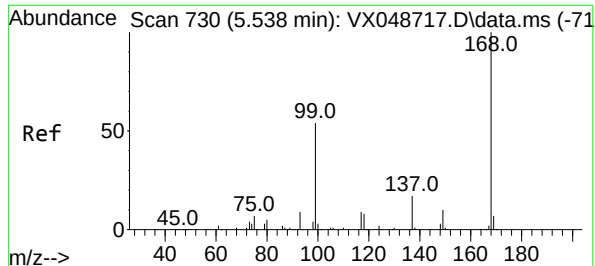
(#) = qualifier out of range (m) = manual integration (+) = signals summed

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120525\
Data File : VX048731.D
Acq On : 05 Dec 2025 13:32
Operator : JC/MD
Sample : Q3757-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-14A-17.5-120325

Quant Time: Dec 05 22:03:35 2025
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X120425W.M
Quant Title : SW846 8260
QLast Update : Fri Dec 05 03:27:34 2025
Response via : Initial Calibration

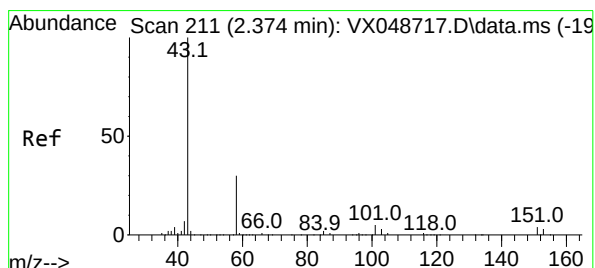
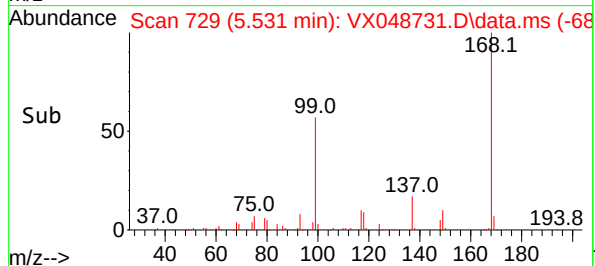
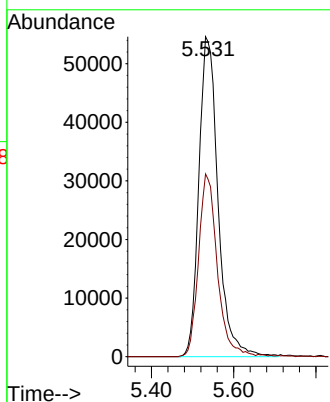
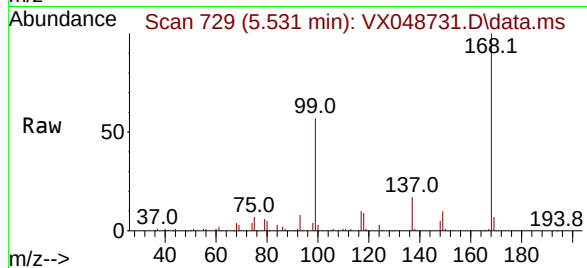




#1
Pentafluorobenzene
Concen: 50.000 ug/l
RT: 5.531 min Scan# 71
Delta R.T. -0.007 min
Lab File: VX048731.D
Acq: 05 Dec 2025 13:32

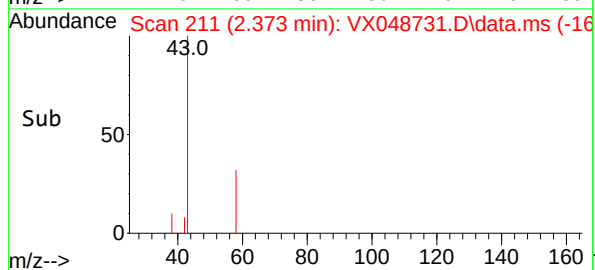
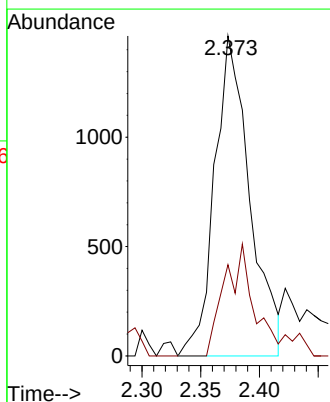
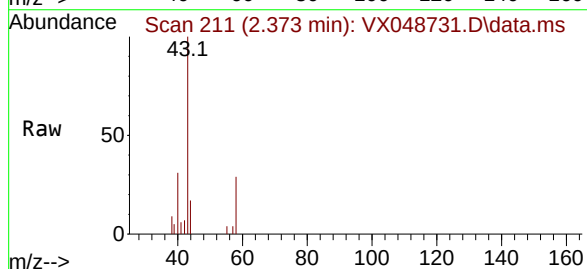
Instrument :
MSVOA_X
ClientSampleId :
MW-14A-17.5-120325

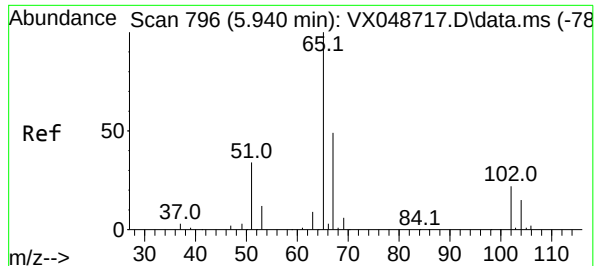
Tgt Ion:168 Resp: 176619
Ion Ratio Lower Upper
168 100
99 57.1 44.2 66.4



#16
Acetone
Concen: 3.368 ug/l
RT: 2.373 min Scan# 211
Delta R.T. -0.000 min
Lab File: VX048731.D
Acq: 05 Dec 2025 13:32

Tgt Ion: 43 Resp: 3056
Ion Ratio Lower Upper
43 100
58 28.5 25.0 37.4





#33

1,2-Dichloroethane-d4

Concen: 52.383 ug/l

RT: 5.940 min Scan# 796

Delta R.T. -0.000 min

Lab File: VX048731.D

Acq: 05 Dec 2025 13:32

Instrument :

MSVOA_X

ClientSampleId :

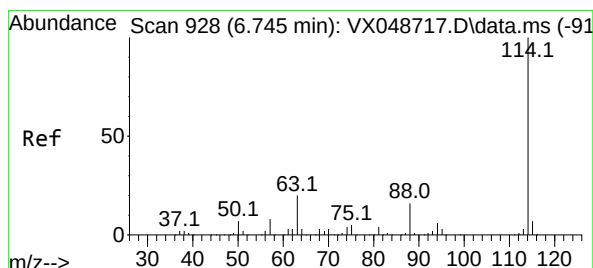
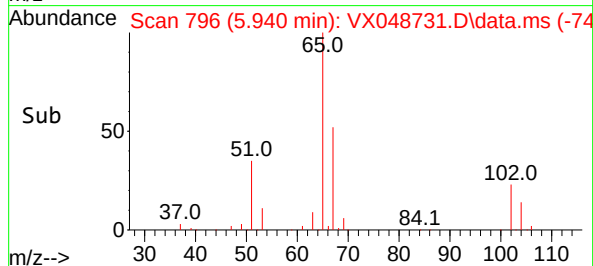
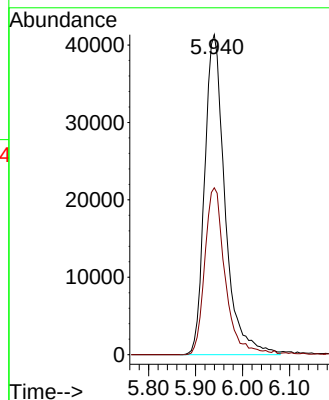
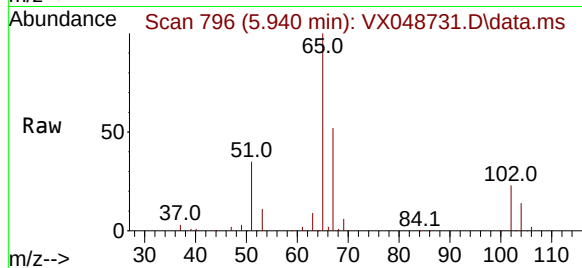
MW-14A-17.5-120325

Tgt Ion: 65 Resp: 124177

Ion Ratio Lower Upper

65 100

67 52.1 0.0 107.4



#34

1,4-Difluorobenzene

Concen: 50.000 ug/l

RT: 6.744 min Scan# 928

Delta R.T. -0.000 min

Lab File: VX048731.D

Acq: 05 Dec 2025 13:32

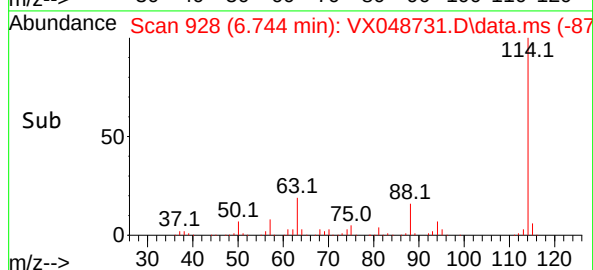
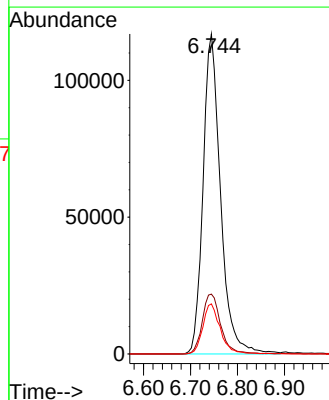
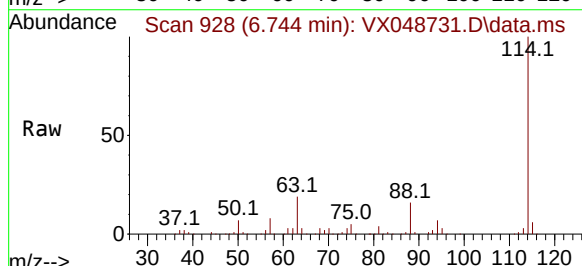
Tgt Ion: 114 Resp: 302604

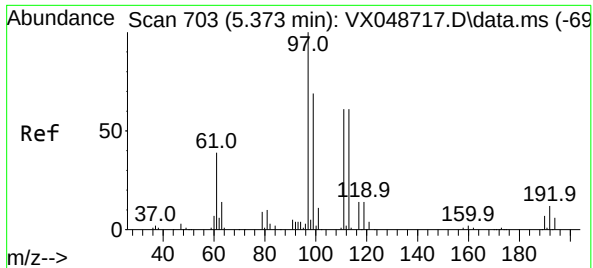
Ion Ratio Lower Upper

114 100

63 18.7 0.0 39.0

88 15.7 0.0 31.8





#35

Dibromofluoromethane

Concen: 49.503 ug/l

RT: 5.367 min Scan# 703

Delta R.T. -0.006 min

Lab File: VX048731.D

Acq: 05 Dec 2025 13:32

Instrument :

MSVOA_X

ClientSampleId :

MW-14A-17.5-120325

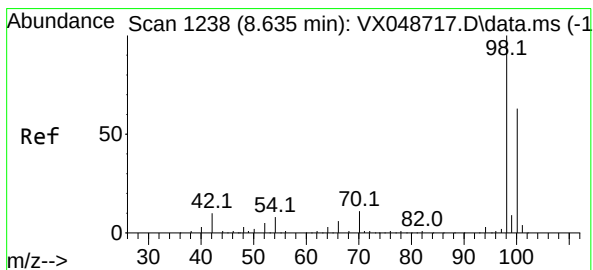
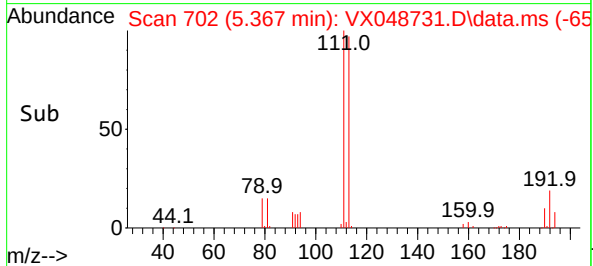
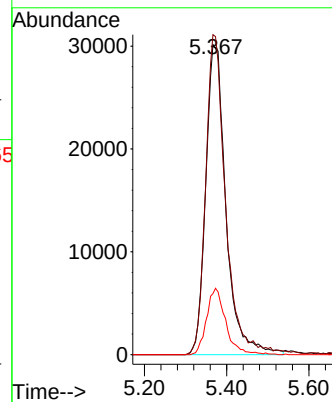
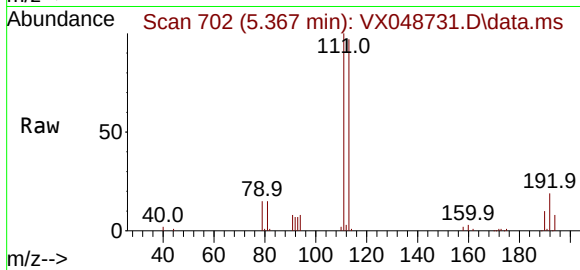
Tgt Ion:113 Resp: 103142

Ion Ratio Lower Upper

113 100

111 102.1 79.5 119.3

192 20.2 16.1 24.1



#50

Toluene-d8

Concen: 48.654 ug/l

RT: 8.628 min Scan# 1237

Delta R.T. -0.006 min

Lab File: VX048731.D

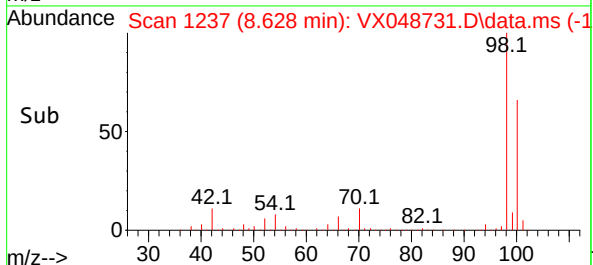
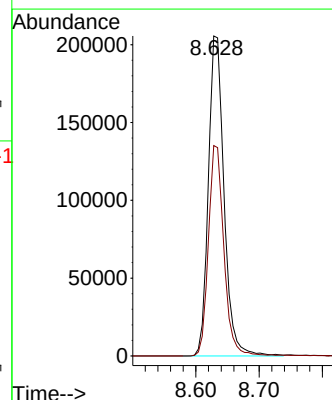
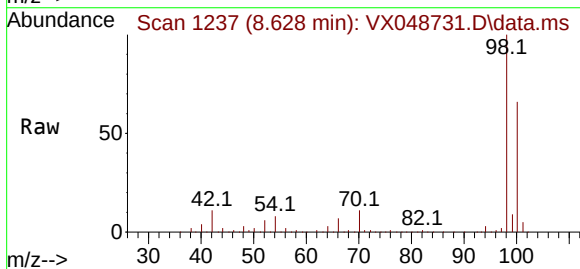
Acq: 05 Dec 2025 13:32

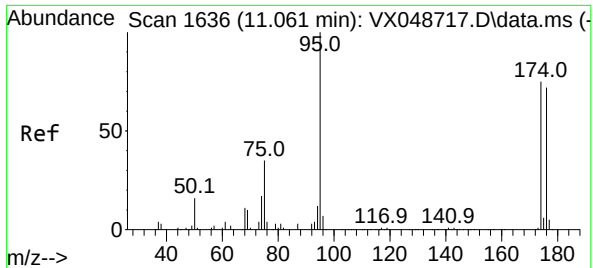
Tgt Ion: 98 Resp: 355343

Ion Ratio Lower Upper

98 100

100 65.2 53.4 80.0

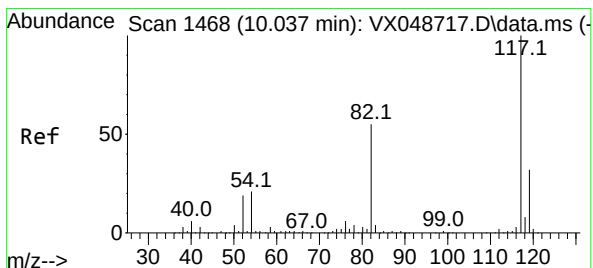
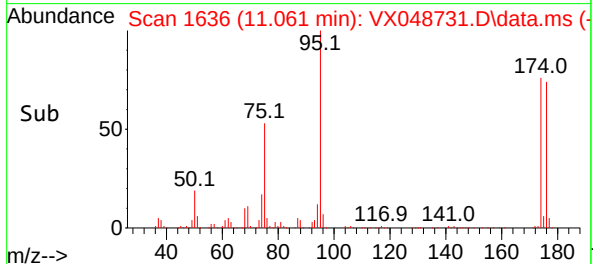
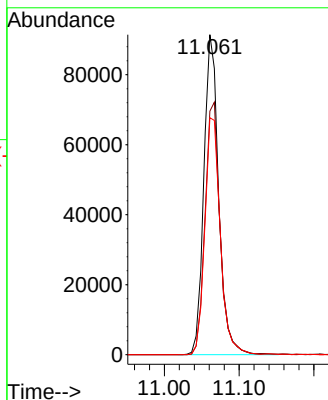
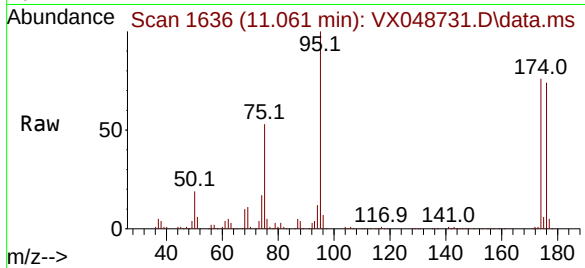




#62
4-Bromofluorobenzene
Concen: 49.979 ug/l
RT: 11.061 min Scan# 1
Delta R.T. -0.000 min
Lab File: VX048731.D
Acq: 05 Dec 2025 13:32

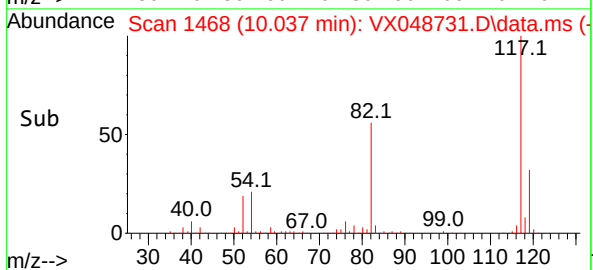
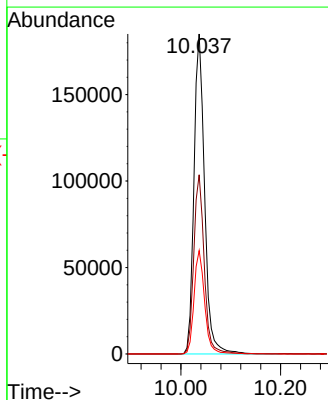
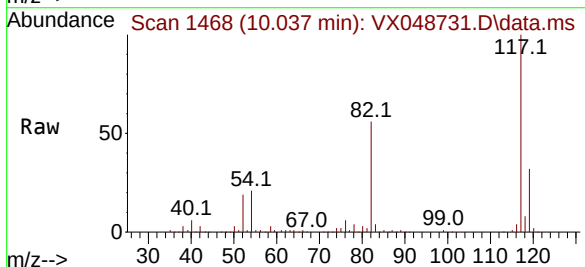
Instrument :
MSVOA_X
ClientSampleId :
MW-14A-17.5-120325

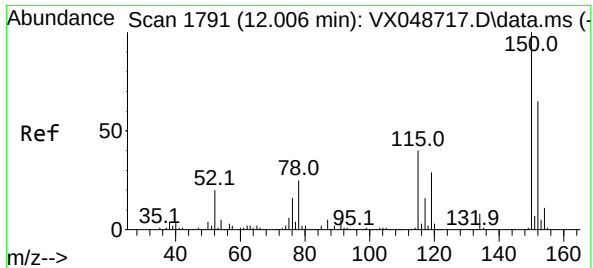
Tgt Ion: 95 Resp: 125871
Ion Ratio Lower Upper
95 100
174 80.3 0.0 157.8
176 77.3 0.0 154.0



#63
Chlorobenzene-d5
Concen: 50.000 ug/l
RT: 10.037 min Scan# 1468
Delta R.T. -0.000 min
Lab File: VX048731.D
Acq: 05 Dec 2025 13:32

Tgt Ion: 117 Resp: 274042
Ion Ratio Lower Upper
117 100
82 56.0 44.1 66.1
119 32.3 25.2 37.8





#72

1,4-Dichlorobenzene-d4

Concen: 50.000 ug/l

RT: 12.006 min Scan# 11

Delta R.T. -0.000 min

Lab File: VX048731.D

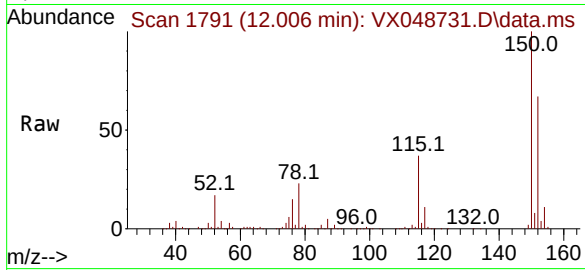
Acq: 05 Dec 2025 13:32

Instrument :

MSVOA_X

ClientSampleId :

MW-14A-17.5-120325



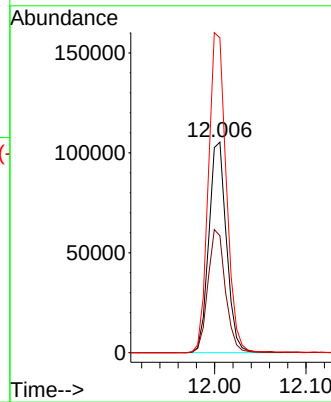
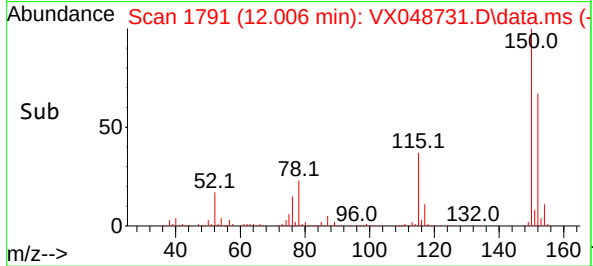
Tgt Ion:152 Resp: 141373

Ion Ratio Lower Upper

152 100

115 58.1 42.1 126.4

150 154.1 0.0 347.8





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

Report of Analysis

Client: JACOBS Engineering Group, Inc.
Project: Former Schlumberger STC PTC Site D3868221
Client Sample ID: BR-03D-490-120325
Lab Sample ID: Q3757-03
Analytical Method: 8260D
Sample Wt/Vol: 5 mL

Level: LOW
Final Vol: 5000 uL

Date Collected: 12/03/25
Date Received: 12/03/25
SDG No.: Q3757
Matrix: Water
% Solid: 0
Test: VOCMS Group3

CAS Number	Parameter	Conc.	Qua.	DF	MDL	LOQ / CRQL	Units	Date Ana.	BatchID
TARGETS									
75-01-4	Vinyl Chloride	3.90		1	0.26	1.00	ug/L	12/05/25 13:53	VX120525
75-35-4	1,1-Dichloroethene	48.1		1	0.23	1.00	ug/L	12/05/25 13:53	VX120525
75-34-3	1,1-Dichloroethane	6.40		1	0.23	1.00	ug/L	12/05/25 13:53	VX120525
156-59-2	cis-1,2-Dichloroethene	34.9		1	0.19	1.00	ug/L	12/05/25 13:53	VX120525
71-55-6	1,1,1-Trichloroethane	0.20	U	1	0.20	1.00	ug/L	12/05/25 13:53	VX120525
71-43-2	Benzene	0.23	J	1	0.15	1.00	ug/L	12/05/25 13:53	VX120525
107-06-2	1,2-Dichloroethane	3.70		1	0.22	1.00	ug/L	12/05/25 13:53	VX120525
79-01-6	Trichloroethene	35.1		1	0.090	1.00	ug/L	12/05/25 13:53	VX120525
79-00-5	1,1,2-Trichloroethane	0.21	U	1	0.21	1.00	ug/L	12/05/25 13:53	VX120525
127-18-4	Tetrachloroethene	0.23	U	1	0.23	1.00	ug/L	12/05/25 13:53	VX120525
SURROGATES									
17060-07-0	1,2-Dichloroethane-d4	55.1			70 (74) - 130 (125)	110%	SPK: 50		
1868-53-7	Dibromofluoromethane	51.8			70 (75) - 130 (124)	104%	SPK: 50		
2037-26-5	Toluene-d8	42.9			70 (86) - 130 (113)	86%	SPK: 50		
460-00-4	4-Bromofluorobenzene	45.5			70 (77) - 130 (121)	91%	SPK: 50		
INTERNAL STANDARDS									
		Area Count							
363-72-4	Pentafluorobenzene	163000							
540-36-3	1,4-Difluorobenzene	281000							
3114-55-4	Chlorobenzene-d5	231000							
3855-82-1	1,4-Dichlorobenzene-d4	110000							

U = Not Detected

LOQ = Limit of Quantitation

MDL = Method Detection Limit

LOD = Limit of Detection

E = Value Exceeds Calibration Range

Q = indicates LCS control criteria did not meet requirements

M = MS/MSD acceptance criteria did not meet requirements

J = Estimated Value

B = Analyte Found in Associated Method Blank

N = Presumptive Evidence of a Compound

* = Values outside of QC limits

D = Dilution

() = Laboratory InHouse Limit

A = Aldol-Condensation Reaction Products

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120525\
 Data File : VX048732.D
 Acq On : 05 Dec 2025 13:53
 Operator : JC/MD
 Sample : Q3757-03
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 BR-03D-490-120325

Quant Time: Dec 05 22:04:09 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X120425W.M
 Quant Title : SW846 8260
 QLast Update : Fri Dec 05 03:27:34 2025
 Response via : Initial Calibration

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

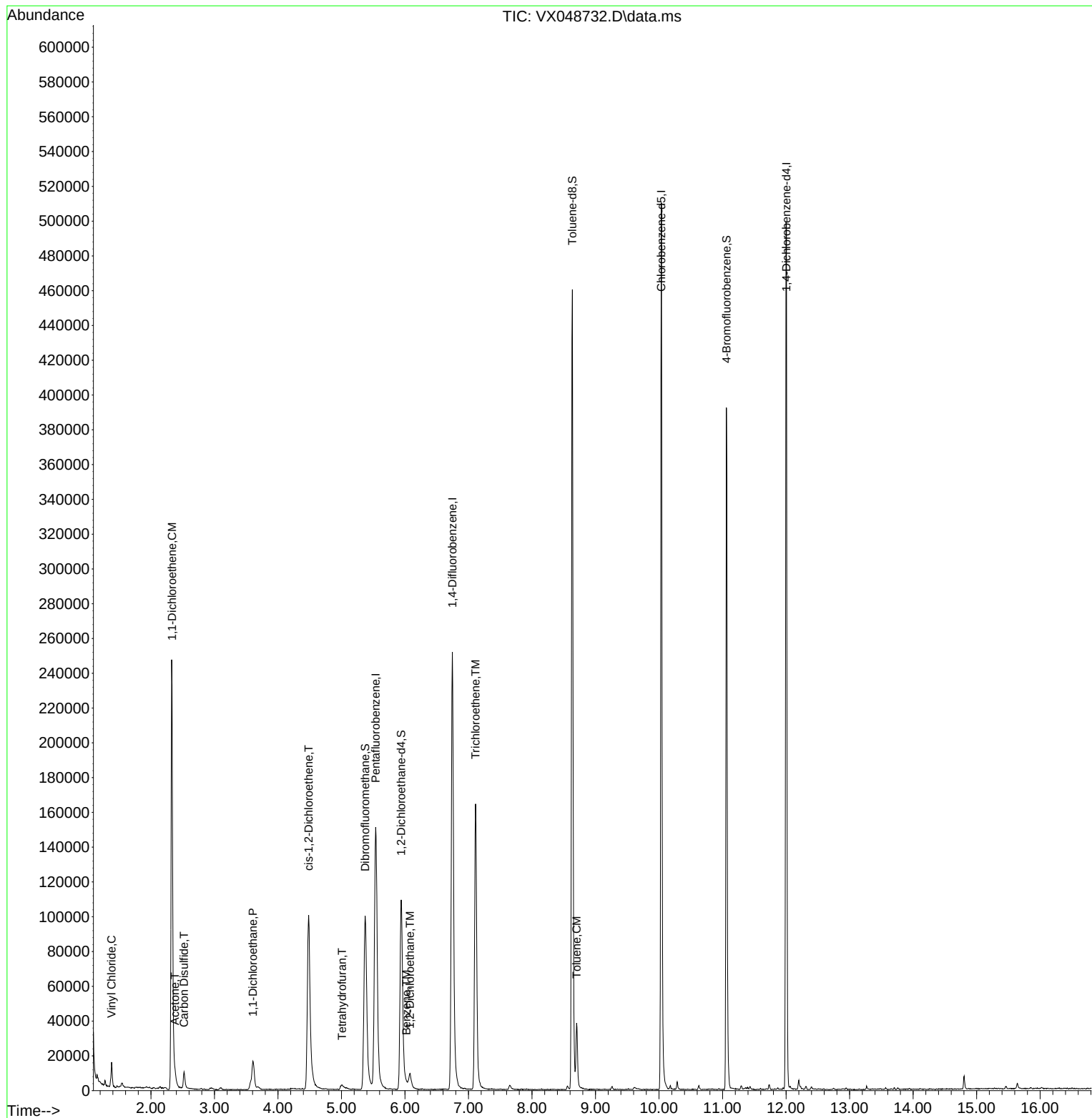
Internal Standards						
1) Pentafluorobenzene	5.538	168	163180	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.745	114	280929	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.037	117	230844	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.006	152	110489	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.940	65	120640	55.082	ug/l	0.00
Spiked Amount	50.000	Range 78 - 117	Recovery	= 110.160%		
35) Dibromofluoromethane	5.373	113	100214	51.809	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	= 103.620%		
50) Toluene-d8	8.635	98	290658	42.868	ug/l	0.00
Spiked Amount	50.000	Range 92 - 112	Recovery	= 85.740%#		
62) 4-Bromofluorobenzene	11.061	95	106460	45.533	ug/l	0.00
Spiked Amount	50.000	Range 83 - 123	Recovery	= 91.060%		
Target Compounds						Qvalue
4) Vinyl Chloride	1.380	62	8761	3.921	ug/l	95
12) 1,1-Dichloroethene	2.331	96	91225	48.143	ug/l	98
16) Acetone	2.380	43	2388	2.848	ug/l	89
17) Carbon Disulfide	2.520	76	14803	2.690	ug/l #	93
24) 1,1-Dichloroethane	3.605	63	21327	6.432	ug/l	99
27) cis-1,2-Dichloroethene	4.483	96	76248	34.939	ug/l	89
29) Tetrahydrofuran	5.007	42	2915	3.085	ug/l #	78
40) Benzene	6.019	78	1766	0.225	ug/l	98
42) 1,2-Dichloroethane	6.080	62	10373	3.732	ug/l	94
44) Trichloroethene	7.110	130	72016	35.143	ug/l	100
52) Toluene	8.702	92	14719	3.108	ug/l	100

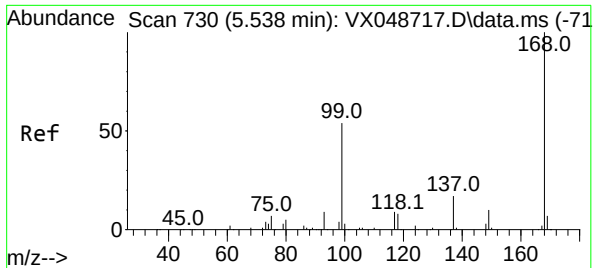
(#) = qualifier out of range (m) = manual integration (+) = signals summed

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120525\
Data File : VX048732.D
Acq On : 05 Dec 2025 13:53
Operator : JC/MD
Sample : Q3757-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
BR-03D-490-120325

Quant Time: Dec 05 22:04:09 2025
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X120425W.M
Quant Title : SW846 8260
QLast Update : Fri Dec 05 03:27:34 2025
Response via : Initial Calibration





#1
Pentafluorobenzene

Concen: 50.000 ug/l

RT: 5.538 min Scan# 71

Delta R.T. -0.000 min

Lab File: VX048732.D

Acq: 05 Dec 2025 13:53

Instrument :

MSVOA_X

ClientSampleId :

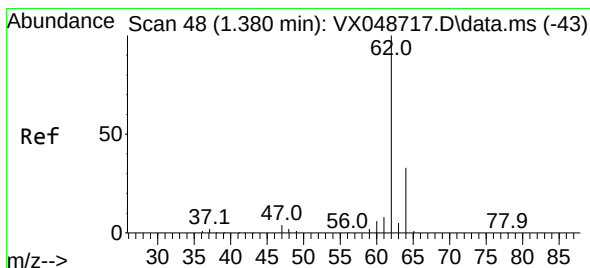
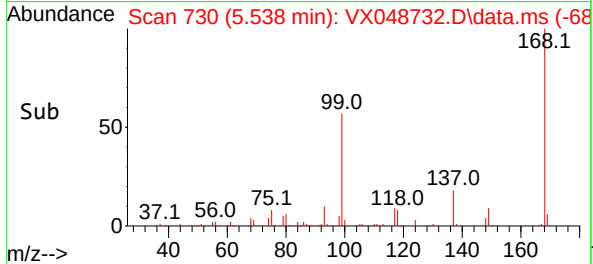
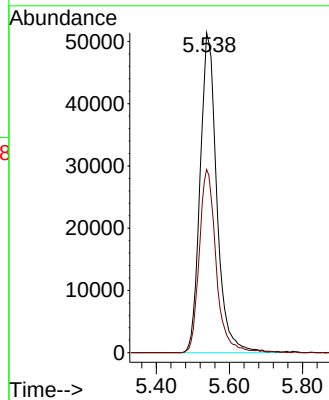
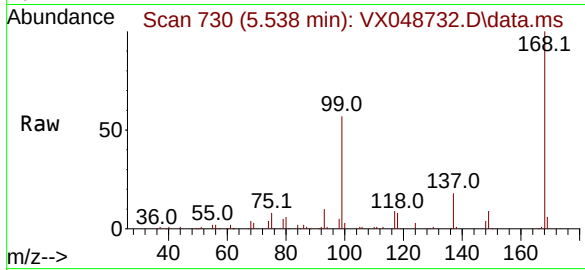
BR-03D-490-120325

Tgt Ion:168 Resp: 163180

Ion Ratio Lower Upper

168 100

99 57.3 44.2 66.4



#4

Vinyl Chloride

Concen: 3.921 ug/l

RT: 1.380 min Scan# 48

Delta R.T. -0.000 min

Lab File: VX048732.D

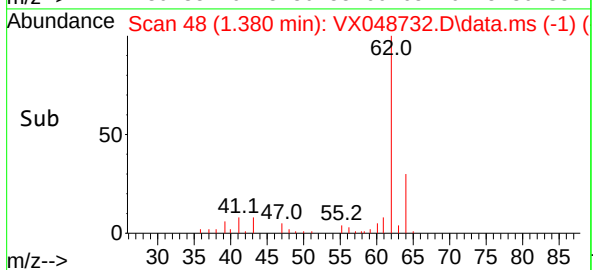
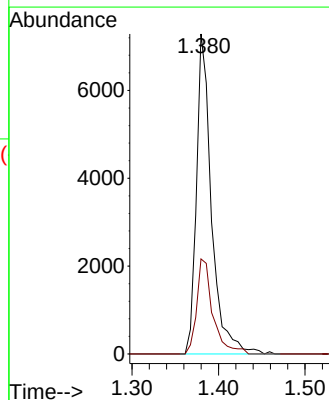
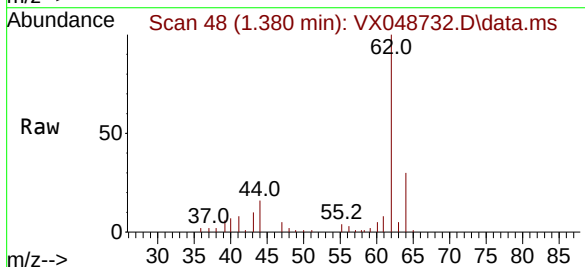
Acq: 05 Dec 2025 13:53

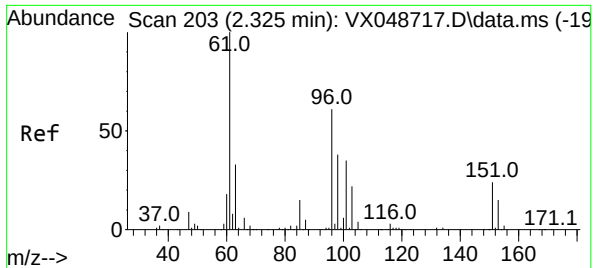
Tgt Ion: 62 Resp: 8761

Ion Ratio Lower Upper

62 100

64 29.7 26.0 39.0





#12

1,1-Dichloroethene

Concen: 48.143 ug/l

RT: 2.331 min Scan# 204

Delta R.T. 0.006 min

Lab File: VX048732.D

Acq: 05 Dec 2025 13:53

Instrument :

MSVOA_X

ClientSampleId :

BR-03D-490-120325

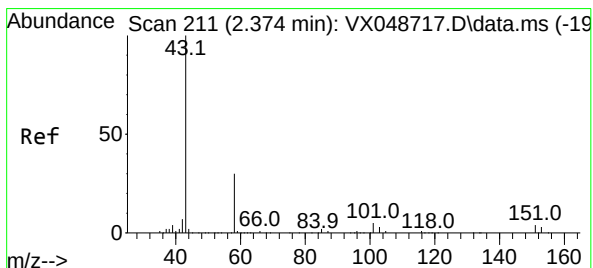
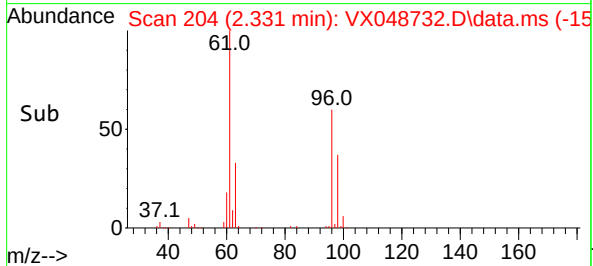
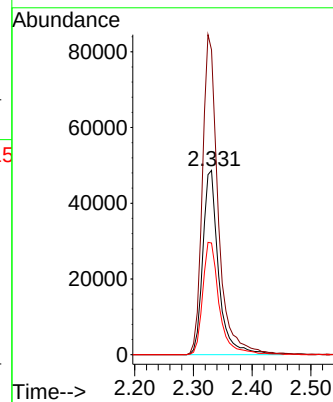
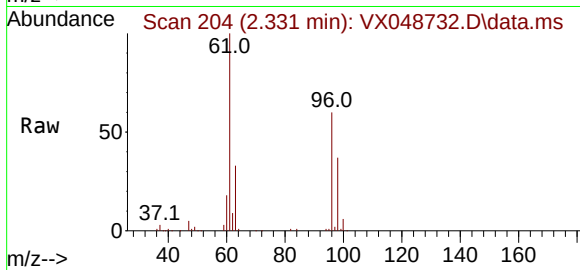
Tgt Ion: 96 Resp: 91225

Ion Ratio Lower Upper

96 100

61 165.4 130.7 196.1

98 60.6 49.8 74.8



#16

Acetone

Concen: 2.848 ug/l

RT: 2.380 min Scan# 212

Delta R.T. 0.006 min

Lab File: VX048732.D

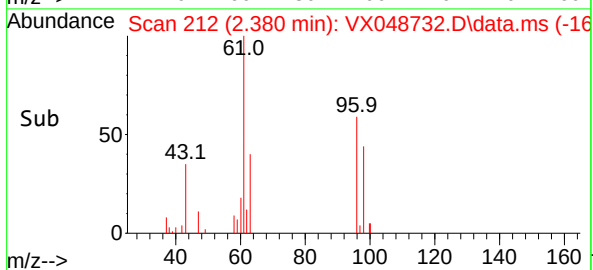
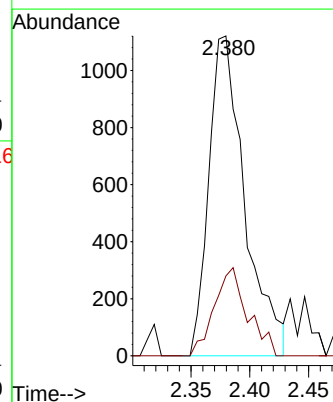
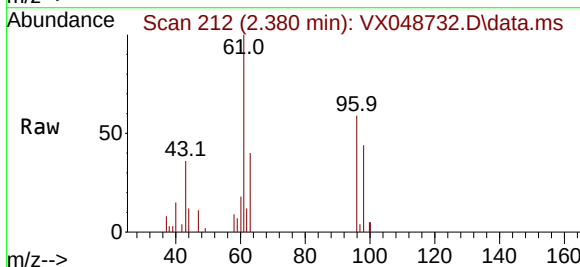
Acq: 05 Dec 2025 13:53

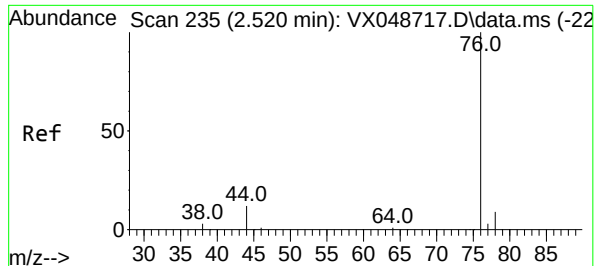
Tgt Ion: 43 Resp: 2388

Ion Ratio Lower Upper

43 100

58 25.0 25.0 37.4

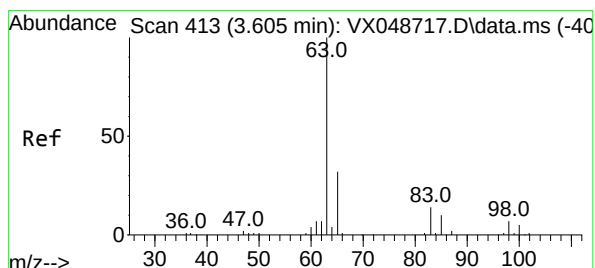
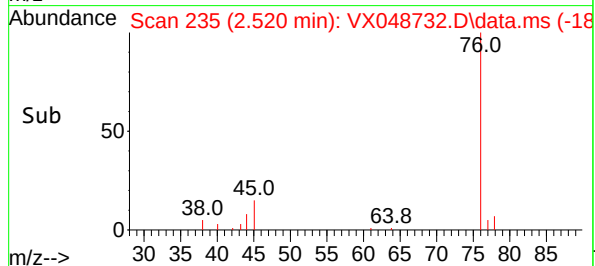
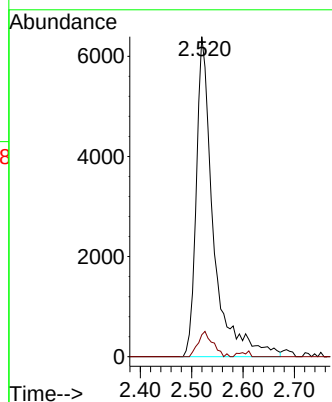
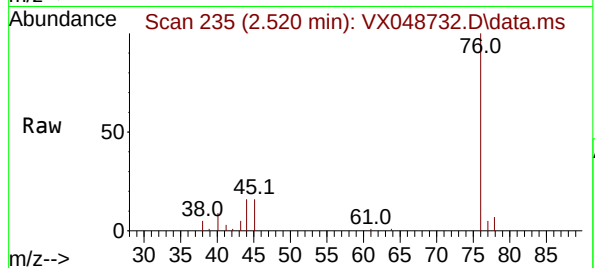




#17
Carbon Disulfide
Concen: 2.690 ug/l
RT: 2.520 min Scan# 21
Delta R.T. -0.000 min
Lab File: VX048732.D
Acq: 05 Dec 2025 13:53

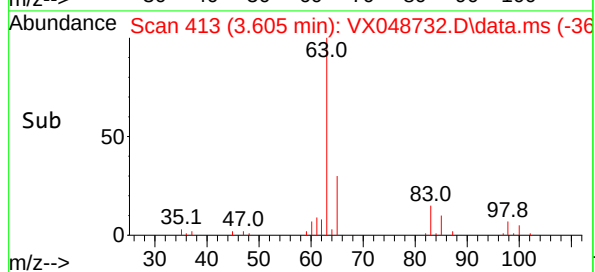
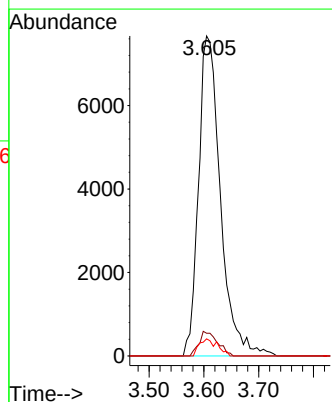
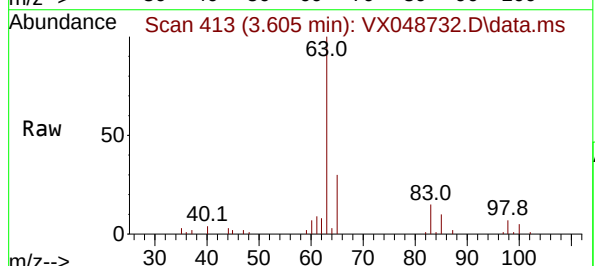
Instrument :
MSVOA_X
ClientSampleId :
BR-03D-490-120325

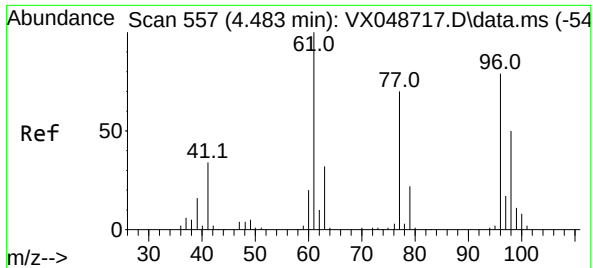
Tgt Ion: 76 Resp: 14803
Ion Ratio Lower Upper
76 100
78 6.8 7.4 11.2#



#24
1,1-Dichloroethane
Concen: 6.432 ug/l
RT: 3.605 min Scan# 413
Delta R.T. -0.000 min
Lab File: VX048732.D
Acq: 05 Dec 2025 13:53

Tgt Ion: 63 Resp: 21327
Ion Ratio Lower Upper
63 100
98 7.1 3.5 10.6
100 5.4 2.5 7.4





#27

cis-1,2-Dichloroethene

Concen: 34.939 ug/l

RT: 4.483 min Scan# 51

Delta R.T. -0.000 min

Lab File: VX048732.D

Acq: 05 Dec 2025 13:53

Instrument :

MSVOA_X

ClientSampleId :

BR-03D-490-120325

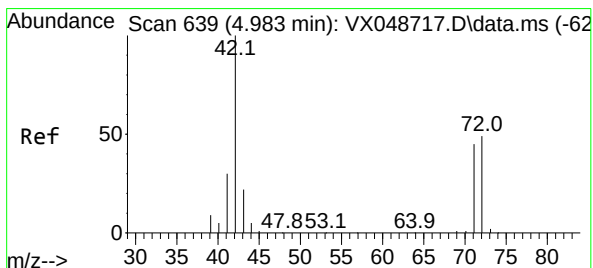
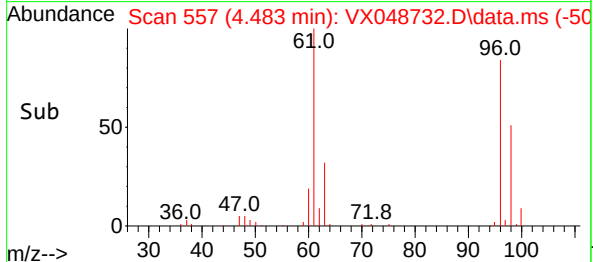
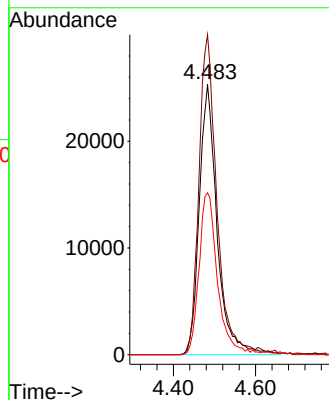
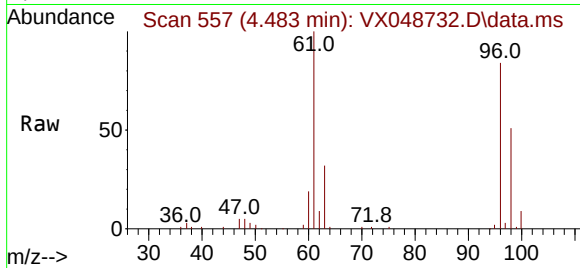
Tgt Ion: 96 Resp: 76248

Ion Ratio Lower Upper

96 100

61 118.2 0.0 271.0

98 62.0 0.0 130.2



#29

Tetrahydrofuran

Concen: 3.085 ug/l

RT: 5.007 min Scan# 643

Delta R.T. 0.024 min

Lab File: VX048732.D

Acq: 05 Dec 2025 13:53

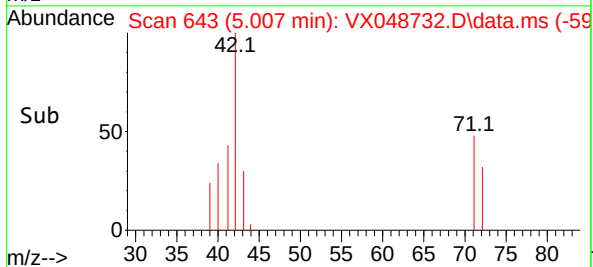
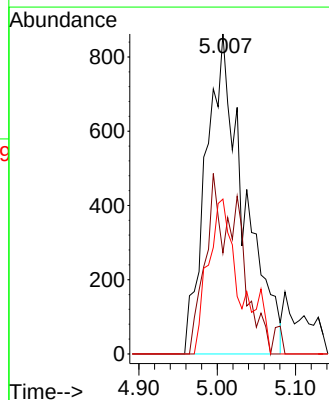
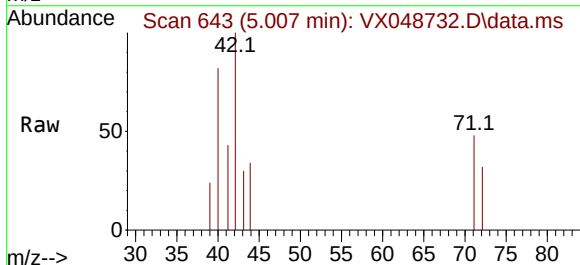
Tgt Ion: 42 Resp: 2915

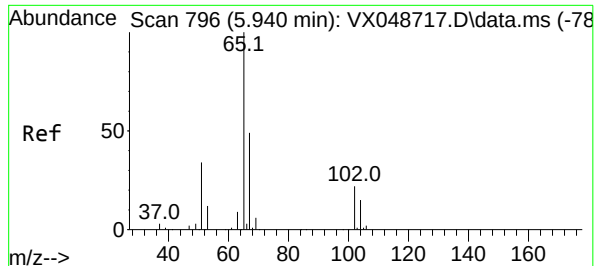
Ion Ratio Lower Upper

42 100

72 24.0 39.4 59.2#

71 40.5 35.8 53.6





#33

1,2-Dichloroethane-d4

Concen: 55.082 ug/l

RT: 5.940 min Scan# 796

Delta R.T. -0.000 min

Lab File: VX048732.D

Acq: 05 Dec 2025 13:53

Instrument :

MSVOA_X

ClientSampleId :

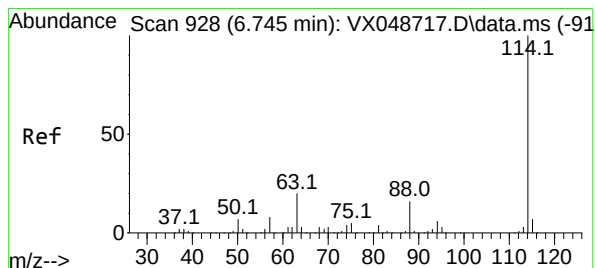
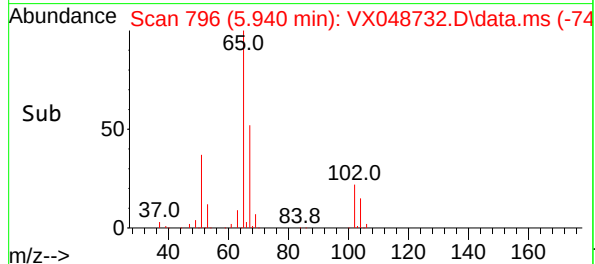
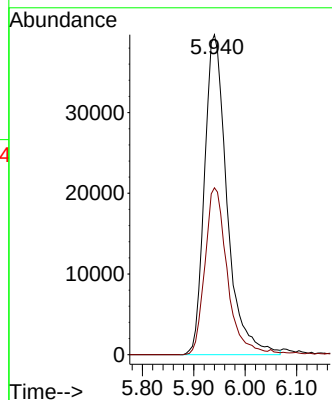
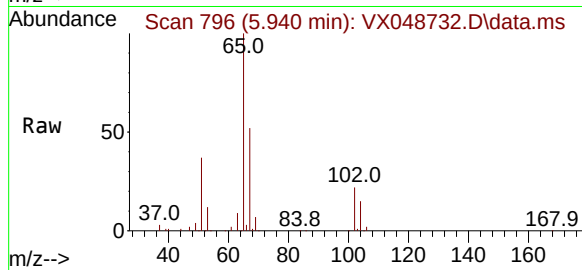
BR-03D-490-120325

Tgt Ion: 65 Resp: 120640

Ion Ratio Lower Upper

65 100

67 51.6 0.0 107.4



#34

1,4-Difluorobenzene

Concen: 50.000 ug/l

RT: 6.745 min Scan# 928

Delta R.T. -0.000 min

Lab File: VX048732.D

Acq: 05 Dec 2025 13:53

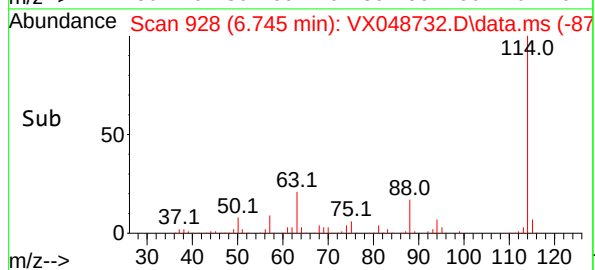
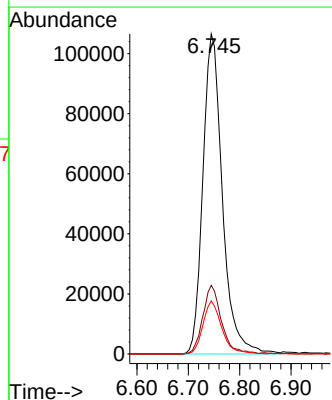
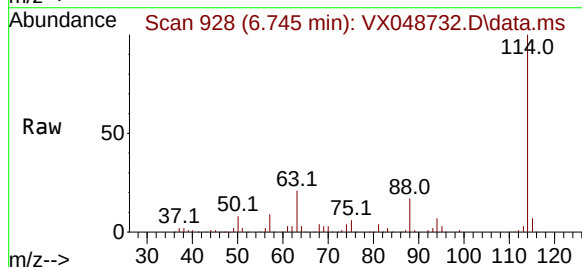
Tgt Ion: 114 Resp: 280929

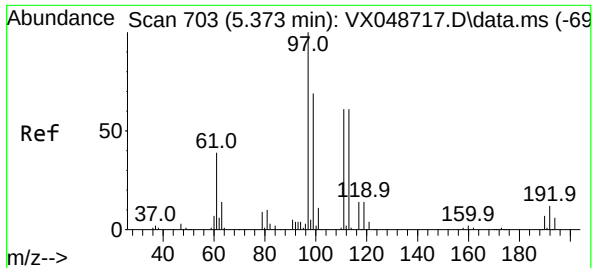
Ion Ratio Lower Upper

114 100

63 21.4 0.0 39.0

88 16.6 0.0 31.8





#35

Dibromofluoromethane

Concen: 51.809 ug/l

RT: 5.373 min Scan# 703

Delta R.T. -0.000 min

Lab File: VX048732.D

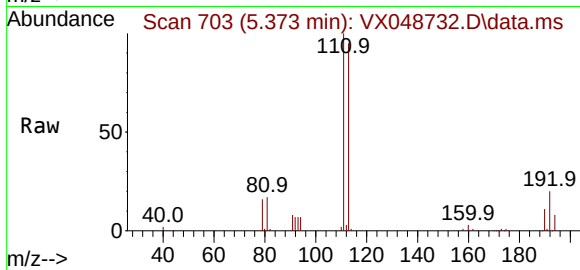
Acq: 05 Dec 2025 13:53

Instrument :

MSVOA_X

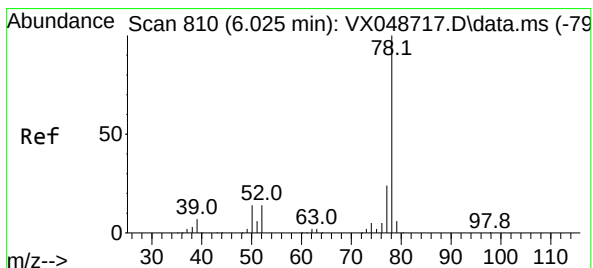
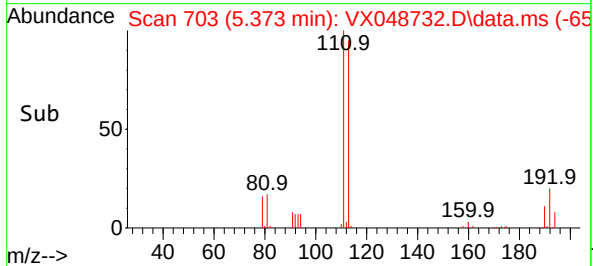
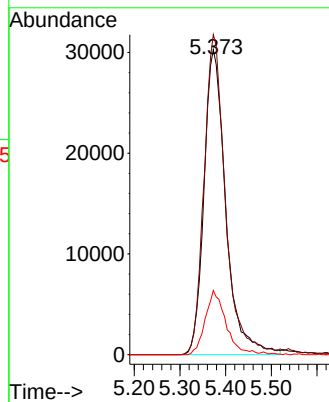
ClientSampleId :

BR-03D-490-120325



Tgt Ion:113 Resp: 100214

Ion	Ratio	Lower	Upper
113	100		
111	102.9	79.5	119.3
192	19.9	16.1	24.1



#40

Benzene

Concen: 0.225 ug/l

RT: 6.019 min Scan# 809

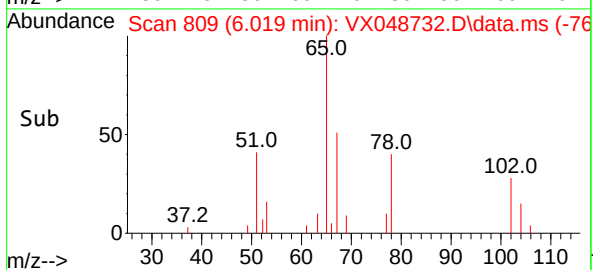
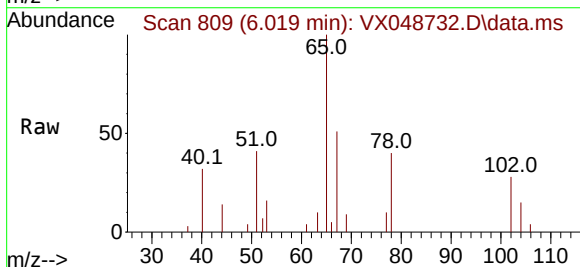
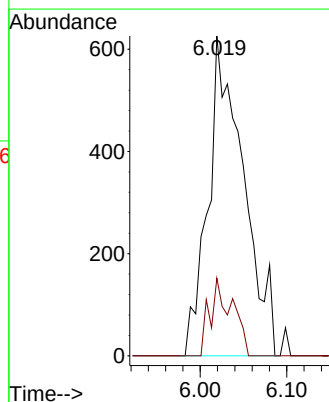
Delta R.T. -0.006 min

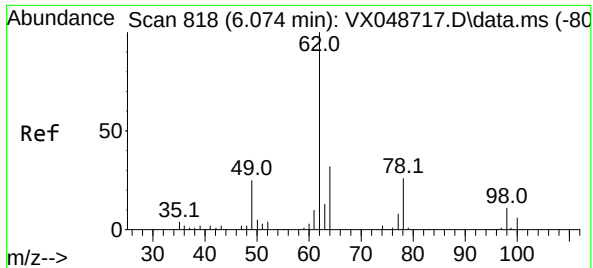
Lab File: VX048732.D

Acq: 05 Dec 2025 13:53

Tgt Ion: 78 Resp: 1766

Ion	Ratio	Lower	Upper
78	100		
77	24.3	18.8	28.2

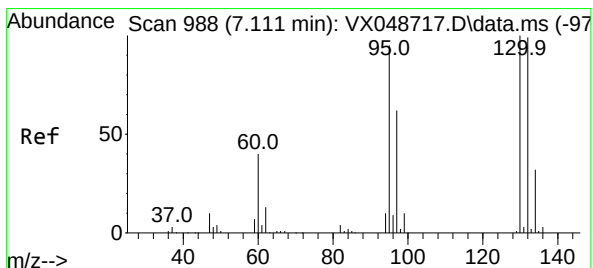
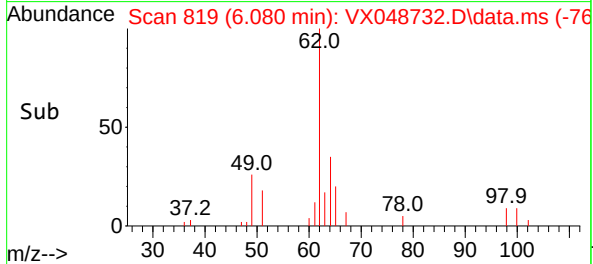
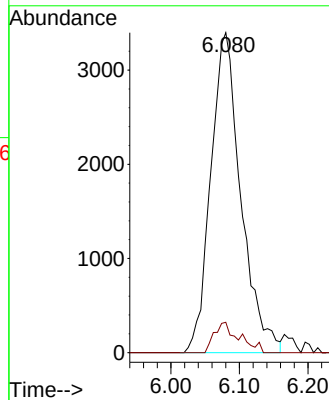
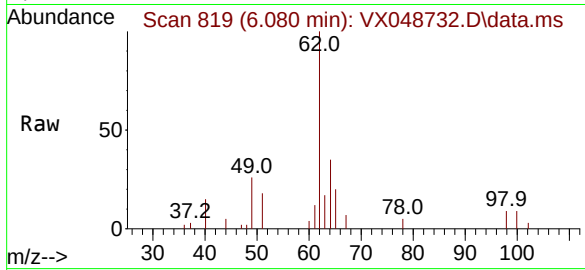




#42
1,2-Dichloroethane
Concen: 3.732 ug/l
RT: 6.080 min Scan# 818
Delta R.T. 0.006 min
Lab File: VX048732.D
Acq: 05 Dec 2025 13:53

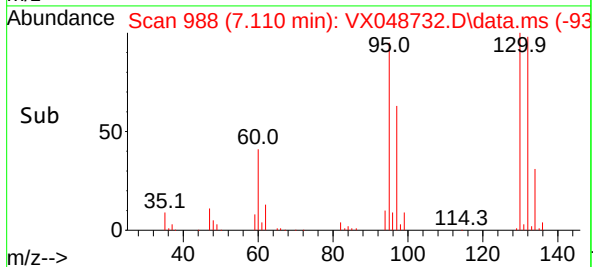
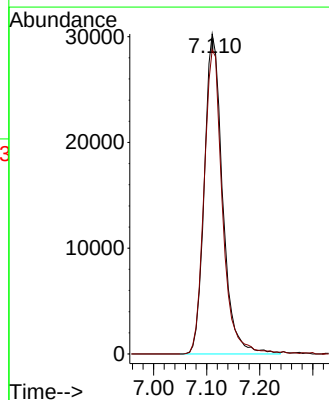
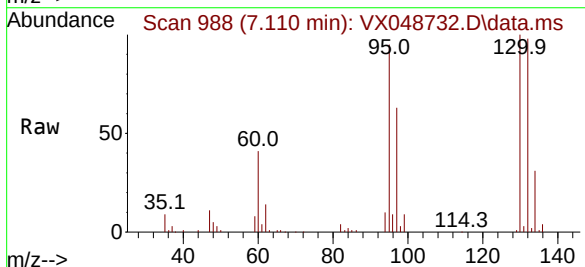
Instrument :
MSVOA_X
ClientSampleId :
BR-03D-490-120325

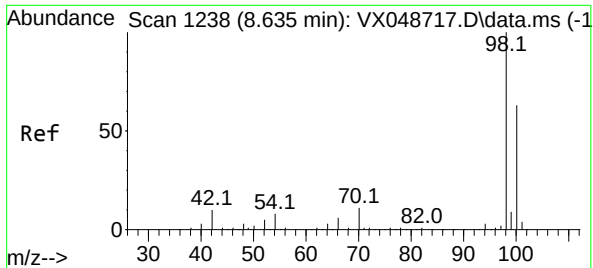
Tgt Ion: 62 Resp: 10373
Ion Ratio Lower Upper
62 100
98 7.8 0.0 20.2



#44
Trichloroethene
Concen: 35.143 ug/l
RT: 7.110 min Scan# 988
Delta R.T. -0.000 min
Lab File: VX048732.D
Acq: 05 Dec 2025 13:53

Tgt Ion:130 Resp: 72016
Ion Ratio Lower Upper
130 100
95 95.0 0.0 189.6





#50

Toluene-d8

Concen: 42.868 ug/l

RT: 8.635 min Scan# 1238

Delta R.T. -0.000 min

Lab File: VX048732.D

Acq: 05 Dec 2025 13:53

Instrument :

MSVOA_X

ClientSampleId :

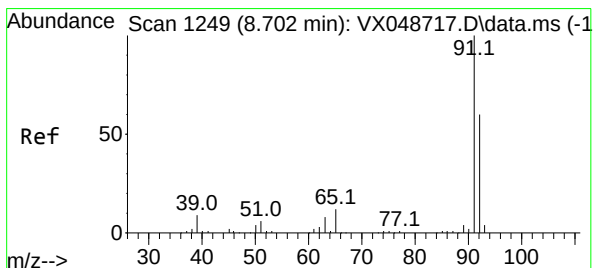
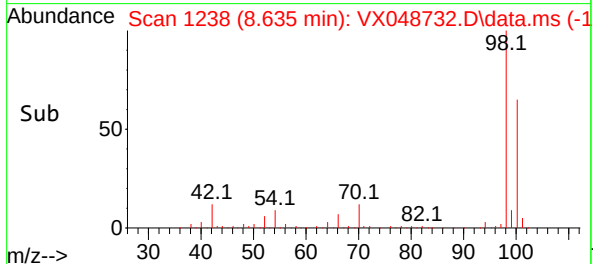
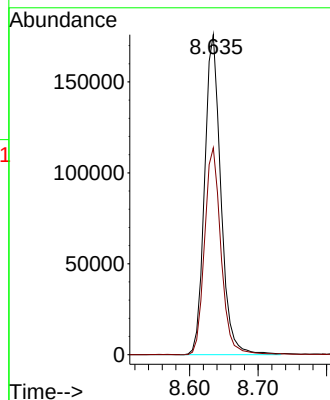
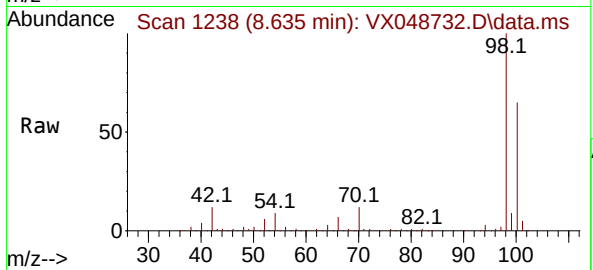
BR-03D-490-120325

Tgt Ion: 98 Resp: 290658

Ion Ratio Lower Upper

98 100

100 65.3 53.4 80.0



#52

Toluene

Concen: 3.108 ug/l

RT: 8.702 min Scan# 1249

Delta R.T. -0.000 min

Lab File: VX048732.D

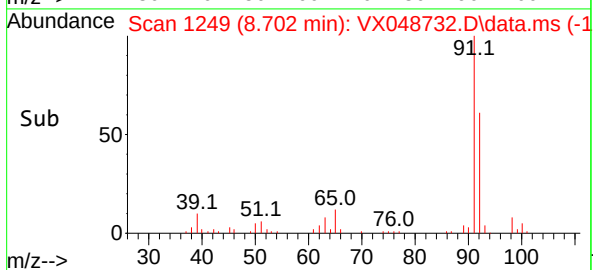
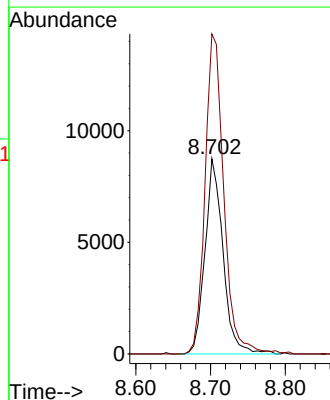
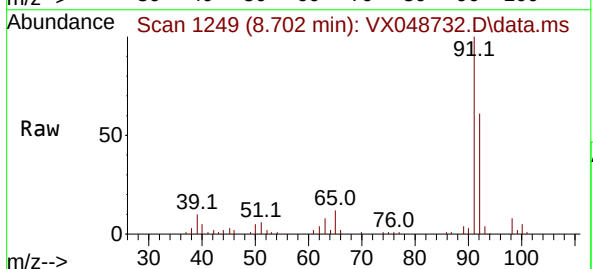
Acq: 05 Dec 2025 13:53

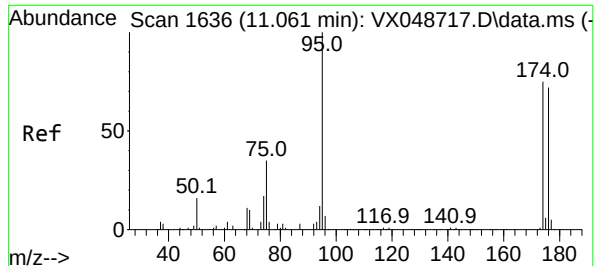
Tgt Ion: 92 Resp: 14719

Ion Ratio Lower Upper

92 100

91 169.9 136.3 204.5





#62

4-Bromofluorobenzene

Concen: 45.533 ug/l

RT: 11.061 min Scan# 1

Delta R.T. -0.000 min

Lab File: VX048732.D

Acq: 05 Dec 2025 13:53

Instrument :

MSVOA_X

ClientSampleId :

BR-03D-490-120325

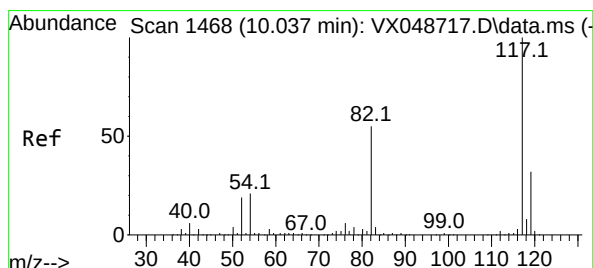
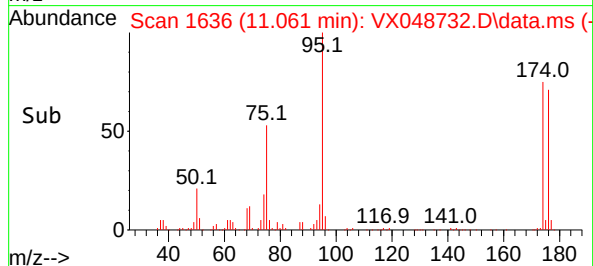
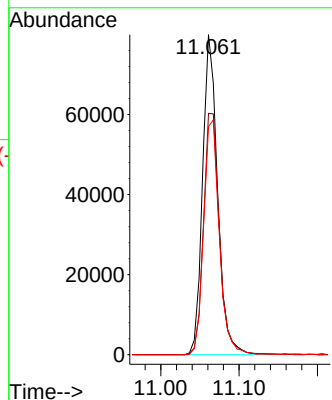
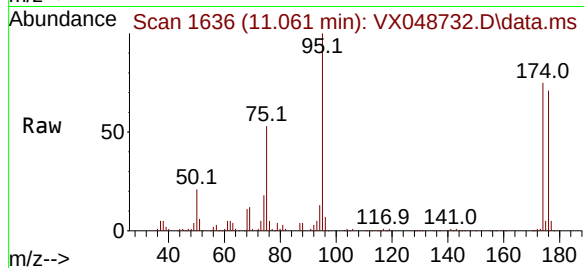
Tgt Ion: 95 Resp: 106460

Ion Ratio Lower Upper

95 100

174 80.3 0.0 157.8

176 76.6 0.0 154.0



#63

Chlorobenzene-d5

Concen: 50.000 ug/l

RT: 10.037 min Scan# 1468

Delta R.T. -0.000 min

Lab File: VX048732.D

Acq: 05 Dec 2025 13:53

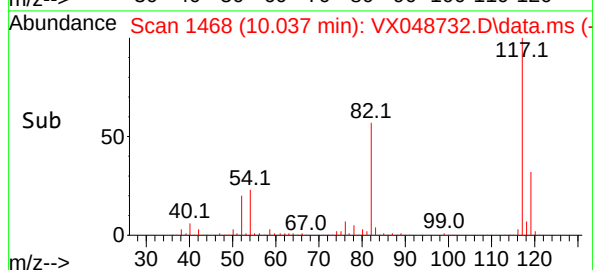
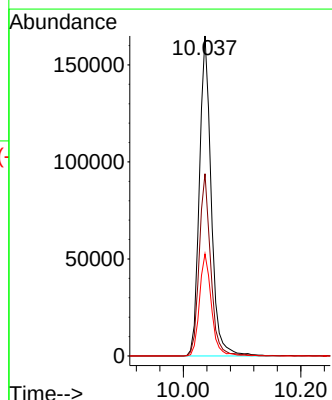
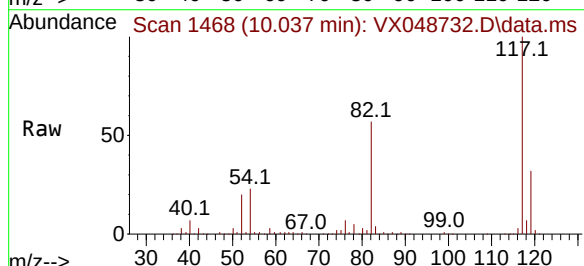
Tgt Ion: 117 Resp: 230844

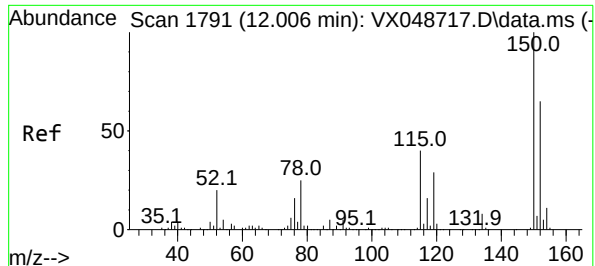
Ion Ratio Lower Upper

117 100

82 56.9 44.1 66.1

119 31.9 25.2 37.8





#72

1,4-Dichlorobenzene-d4

Concen: 50.000 ug/l

RT: 12.006 min Scan# 11

Delta R.T. -0.000 min

Lab File: VX048732.D

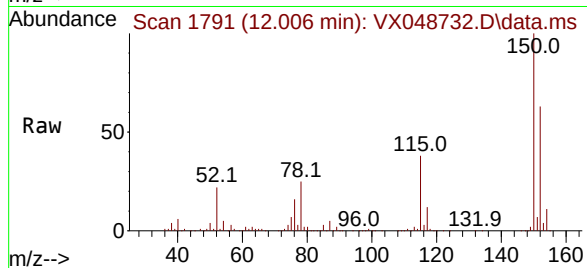
Acq: 05 Dec 2025 13:53

Instrument :

MSVOA_X

ClientSampleId :

BR-03D-490-120325



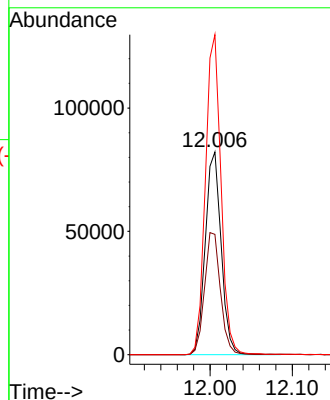
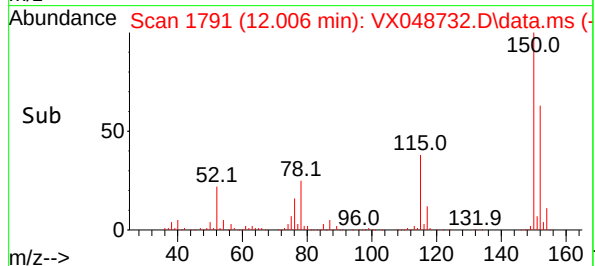
Tgt Ion:152 Resp: 110489

Ion Ratio Lower Upper

152 100

115 60.4 42.1 126.4

150 154.7 0.0 347.8





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

Report of Analysis

Client: JACOBS Engineering Group, Inc.
Project: Former Schlumberger STC PTC Site D3868221
Client Sample ID: MW-14B-46-120325
Lab Sample ID: Q3757-05
Analytical Method: 8260D
Sample Wt/Vol: 5 mL

Level: LOW
Final Vol: 5000 uL

Date Collected: 12/03/25
Date Received: 12/03/25
SDG No.: Q3757
Matrix: Water
% Solid: 0
Test: VOCMS Group3

CAS Number	Parameter	Conc.	Qua.	DF	MDL	LOQ / CRQL	Units	Date Ana.	BatchID
TARGETS									
75-01-4	Vinyl Chloride	0.26	U	1	0.26	1.00	ug/L	12/05/25 14:13	VX120525
75-35-4	1,1-Dichloroethene	0.23	U	1	0.23	1.00	ug/L	12/05/25 14:13	VX120525
75-34-3	1,1-Dichloroethane	0.23	U	1	0.23	1.00	ug/L	12/05/25 14:13	VX120525
156-59-2	cis-1,2-Dichloroethene	0.19	U	1	0.19	1.00	ug/L	12/05/25 14:13	VX120525
71-55-6	1,1,1-Trichloroethane	0.20	U	1	0.20	1.00	ug/L	12/05/25 14:13	VX120525
71-43-2	Benzene	0.15	U	1	0.15	1.00	ug/L	12/05/25 14:13	VX120525
107-06-2	1,2-Dichloroethane	0.22	U	1	0.22	1.00	ug/L	12/05/25 14:13	VX120525
79-01-6	Trichloroethene	0.090	U	1	0.090	1.00	ug/L	12/05/25 14:13	VX120525
79-00-5	1,1,2-Trichloroethane	0.21	U	1	0.21	1.00	ug/L	12/05/25 14:13	VX120525
127-18-4	Tetrachloroethene	0.23	U	1	0.23	1.00	ug/L	12/05/25 14:13	VX120525
SURROGATES									
17060-07-0	1,2-Dichloroethane-d4	47.2			70 (74) - 130 (125)	94%	SPK: 50		
1868-53-7	Dibromofluoromethane	40.9			70 (75) - 130 (124)	82%	SPK: 50		
2037-26-5	Toluene-d8	35.3			70 (86) - 130 (113)	71%	SPK: 50		
460-00-4	4-Bromofluorobenzene	43.5			70 (77) - 130 (121)	87%	SPK: 50		
INTERNAL STANDARDS									
		Area Count							
363-72-4	Pentafluorobenzene	162000							
540-36-3	1,4-Difluorobenzene	287000							
3114-55-4	Chlorobenzene-d5	220000							
3855-82-1	1,4-Dichlorobenzene-d4	131000							

U = Not Detected

LOQ = Limit of Quantitation

MDL = Method Detection Limit

LOD = Limit of Detection

E = Value Exceeds Calibration Range

Q = indicates LCS control criteria did not meet requirements

M = MS/MSD acceptance criteria did not meet requirements

J = Estimated Value

B = Analyte Found in Associated Method Blank

N = Presumptive Evidence of a Compound

* = Values outside of QC limits

D = Dilution

() = Laboratory InHouse Limit

A = Aldol-Condensation Reaction Products

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120525\
 Data File : VX048733.D
 Acq On : 05 Dec 2025 14:13
 Operator : JC/MD
 Sample : Q3757-05
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-14B-46-120325

Quant Time: Dec 05 22:04:43 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X120425W.M
 Quant Title : SW846 8260
 QLast Update : Fri Dec 05 03:27:34 2025
 Response via : Initial Calibration

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

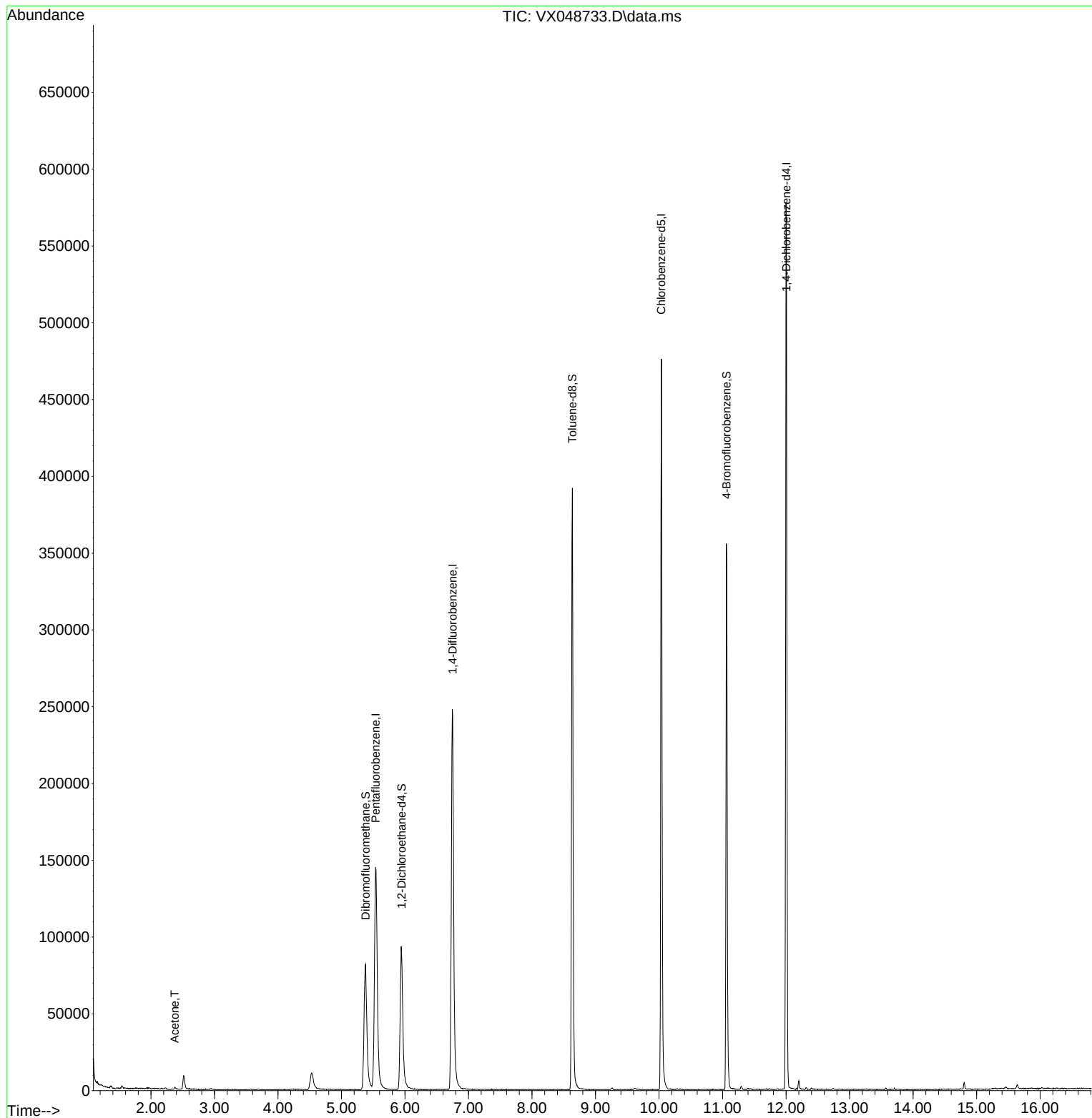
Internal Standards						
1) Pentafluorobenzene	5.537	168	161591	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.751	114	286549	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.037	117	220300	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.006	152	131055	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.946	65	102428	47.227	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	94.460%
35) Dibromofluoromethane	5.379	113	80744	40.924	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	81.840%
50) Toluene-d8	8.634	98	244075	35.291	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	70.580%#
62) 4-Bromofluorobenzene	11.061	95	103850	43.546	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	87.100%
Target Compounds						
16) Acetone	2.373	43	2569	3.094	ug/l	Qvalue 100

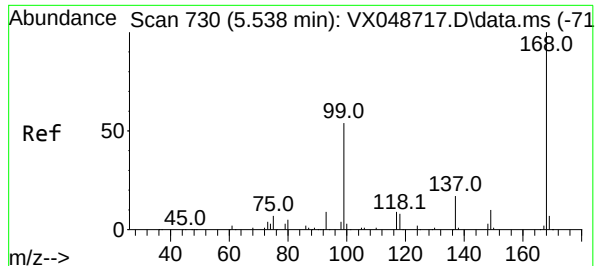
(#) = qualifier out of range (m) = manual integration (+) = signals summed

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120525\
Data File : VX048733.D
Acq On : 05 Dec 2025 14:13
Operator : JC/MD
Sample : Q3757-05
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-14B-46-120325

Quant Time: Dec 05 22:04:43 2025
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X120425W.M
Quant Title : SW846 8260
QLast Update : Fri Dec 05 03:27:34 2025
Response via : Initial Calibration

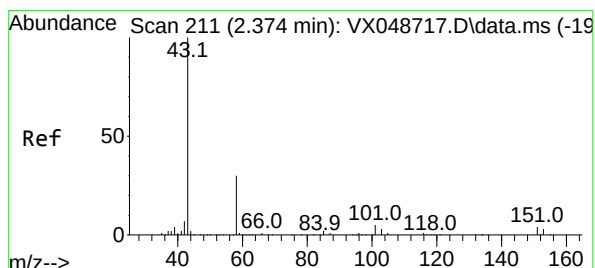
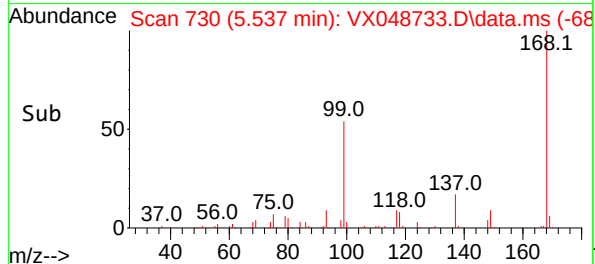
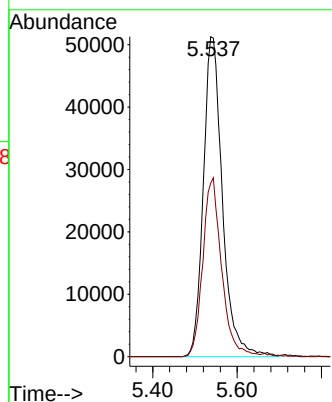
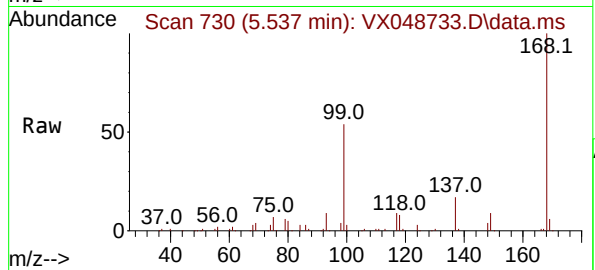




#1
Pentafluorobenzene
Concen: 50.000 ug/l
RT: 5.537 min Scan# 71
Delta R.T. -0.001 min
Lab File: VX048733.D
Acq: 05 Dec 2025 14:13

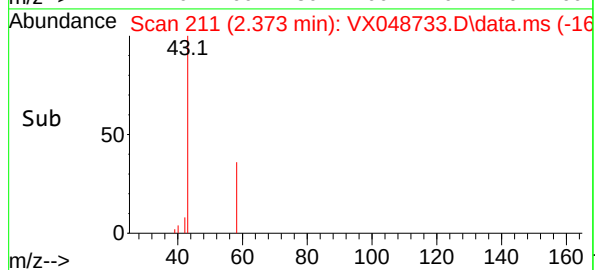
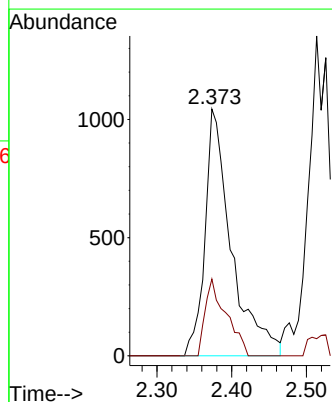
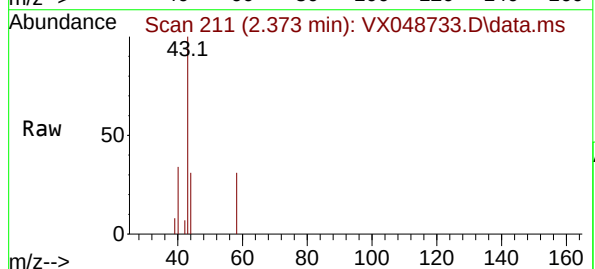
Instrument :
MSVOA_X
ClientSampleId :
MW-14B-46-120325

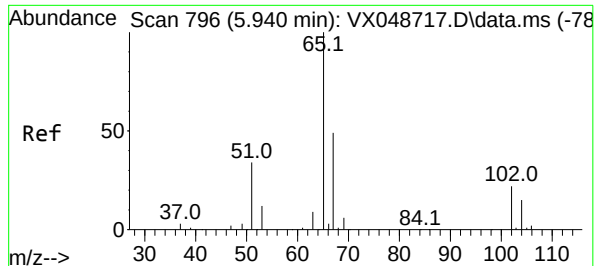
Tgt Ion:168 Resp: 161591
Ion Ratio Lower Upper
168 100
99 54.2 44.2 66.4



#16
Acetone
Concen: 3.094 ug/l
RT: 2.373 min Scan# 211
Delta R.T. -0.000 min
Lab File: VX048733.D
Acq: 05 Dec 2025 14:13

Tgt Ion: 43 Resp: 2569
Ion Ratio Lower Upper
43 100
58 31.1 25.0 37.4





#33

1,2-Dichloroethane-d4

Concen: 47.227 ug/l

RT: 5.946 min Scan# 796

Delta R.T. 0.006 min

Lab File: VX048733.D

Acq: 05 Dec 2025 14:13

Instrument :

MSVOA_X

ClientSampleId :

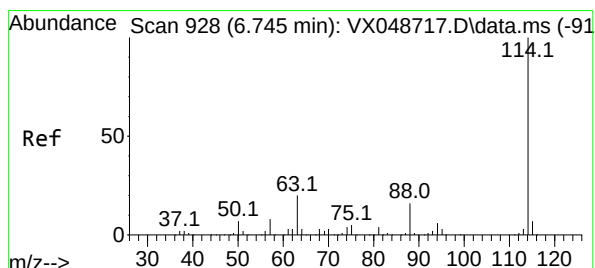
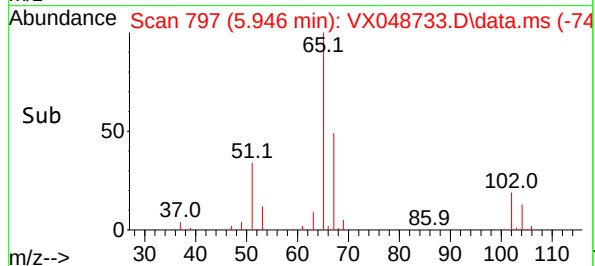
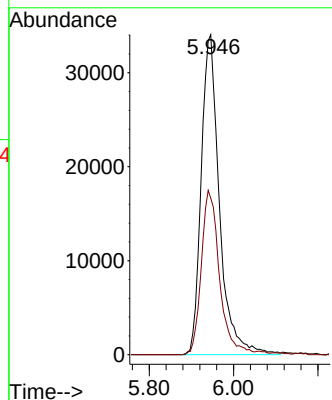
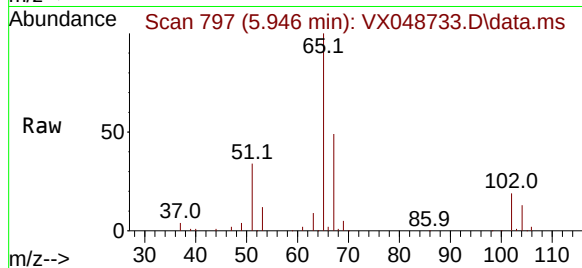
MW-14B-46-120325

Tgt Ion: 65 Resp: 102428

Ion Ratio Lower Upper

65 100

67 51.2 0.0 107.4



#34

1,4-Difluorobenzene

Concen: 50.000 ug/l

RT: 6.751 min Scan# 929

Delta R.T. 0.006 min

Lab File: VX048733.D

Acq: 05 Dec 2025 14:13

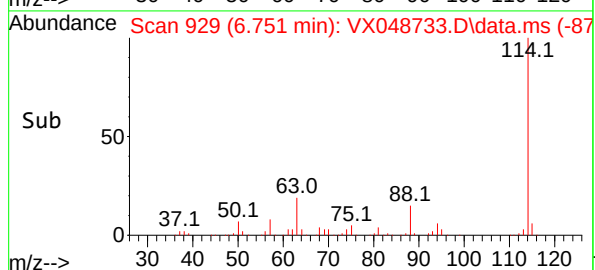
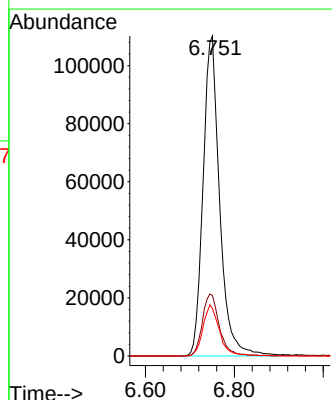
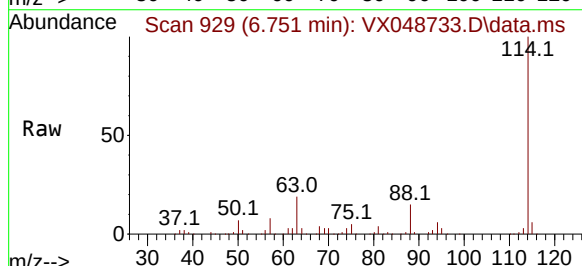
Tgt Ion: 114 Resp: 286549

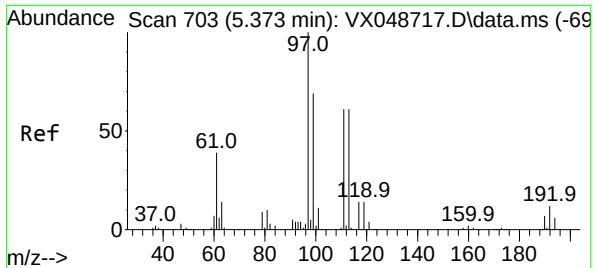
Ion Ratio Lower Upper

114 100

63 18.9 0.0 39.0

88 14.9 0.0 31.8





#35

Dibromofluoromethane

Concen: 40.924 ug/l

RT: 5.379 min Scan# 703

Delta R.T. 0.006 min

Lab File: VX048733.D

Acq: 05 Dec 2025 14:13

Instrument :

MSVOA_X

ClientSampleId :

MW-14B-46-120325

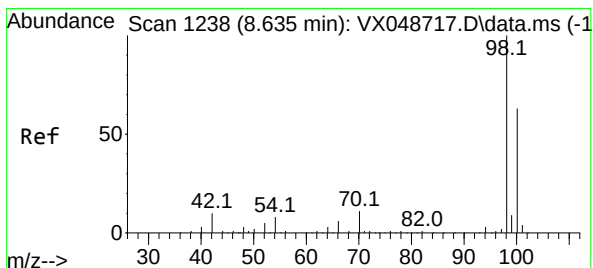
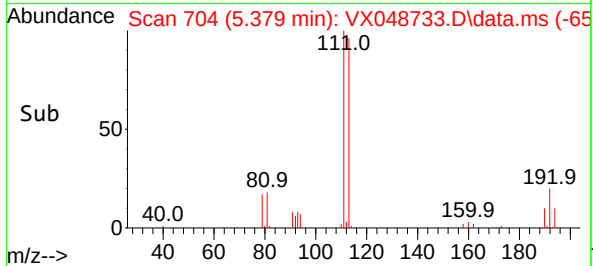
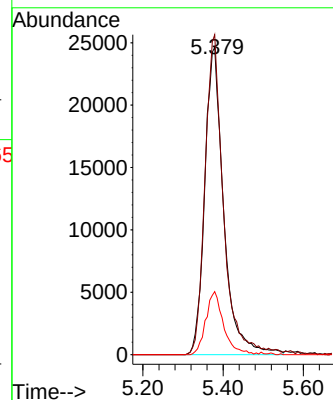
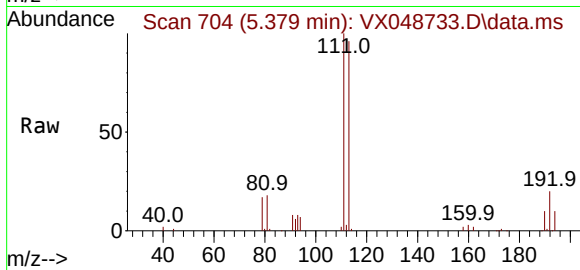
Tgt Ion:113 Resp: 80744

Ion Ratio Lower Upper

113 100

111 101.6 79.5 119.3

192 19.3 16.1 24.1



#50

Toluene-d8

Concen: 35.291 ug/l

RT: 8.634 min Scan# 1238

Delta R.T. -0.000 min

Lab File: VX048733.D

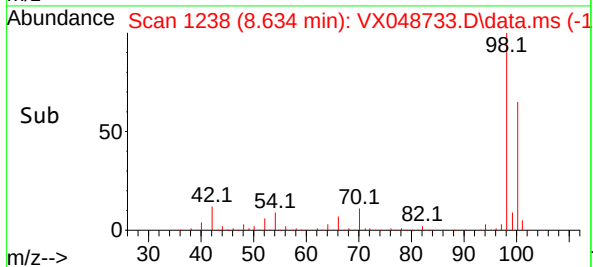
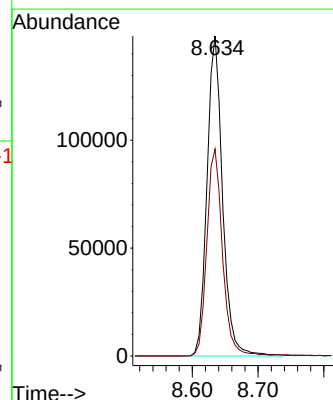
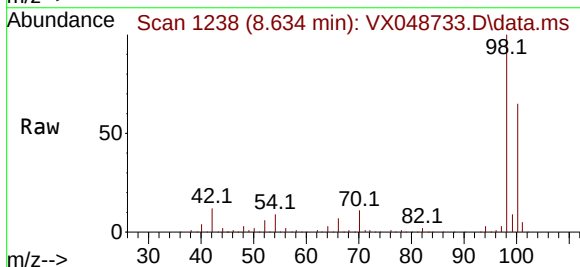
Acq: 05 Dec 2025 14:13

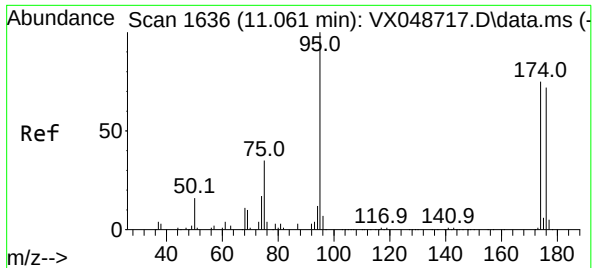
Tgt Ion: 98 Resp: 244075

Ion Ratio Lower Upper

98 100

100 64.9 53.4 80.0

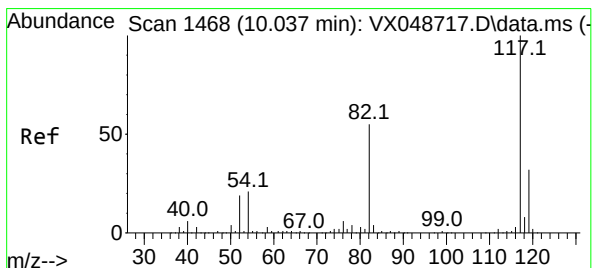
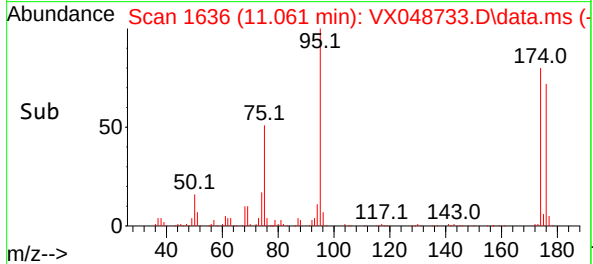
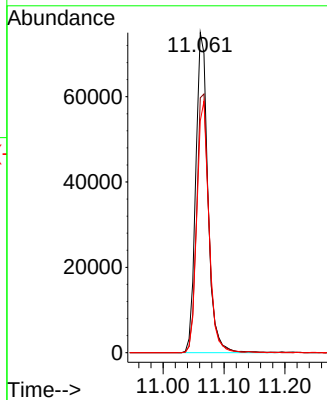
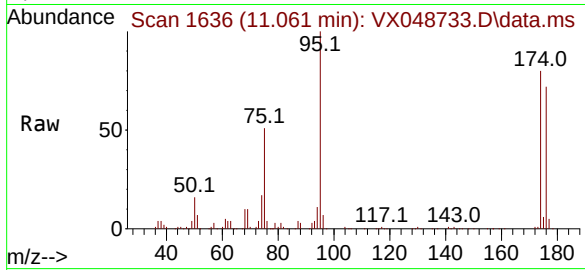




#62
4-Bromofluorobenzene
Concen: 43.546 ug/l
RT: 11.061 min Scan# 1
Delta R.T. -0.000 min
Lab File: VX048733.D
Acq: 05 Dec 2025 14:13

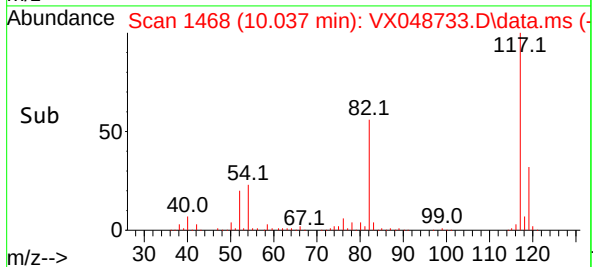
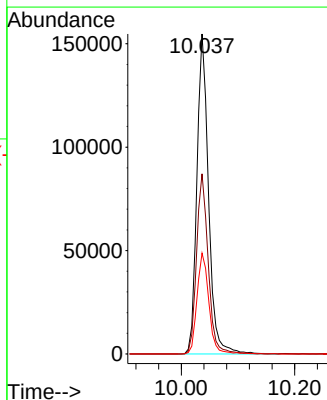
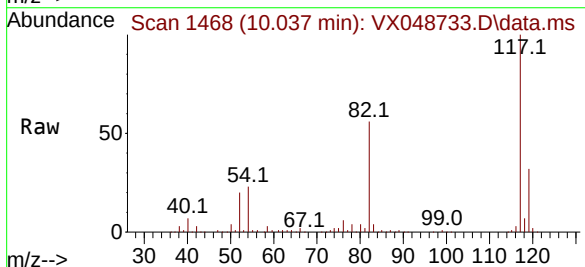
Instrument :
MSVOA_X
ClientSampleId :
MW-14B-46-120325

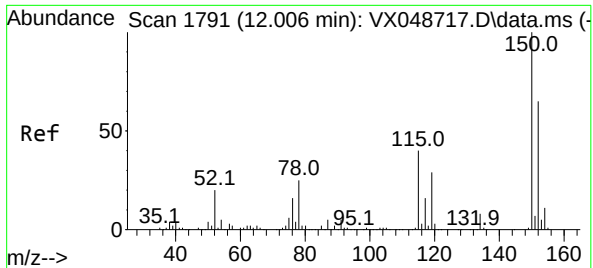
Tgt Ion: 95 Resp: 103850
Ion Ratio Lower Upper
95 100
174 81.4 0.0 157.8
176 78.1 0.0 154.0



#63
Chlorobenzene-d5
Concen: 50.000 ug/l
RT: 10.037 min Scan# 1468
Delta R.T. -0.000 min
Lab File: VX048733.D
Acq: 05 Dec 2025 14:13

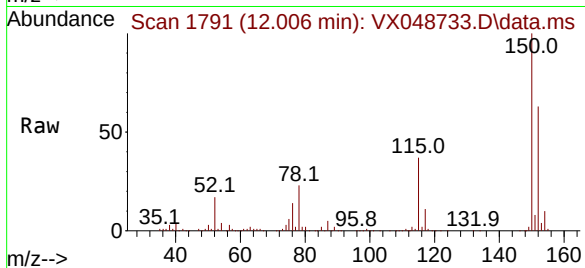
Tgt Ion: 117 Resp: 220300
Ion Ratio Lower Upper
117 100
82 56.3 44.1 66.1
119 31.5 25.2 37.8



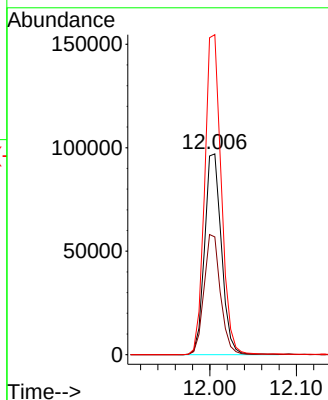
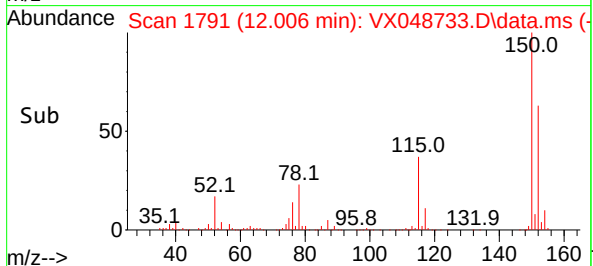


#72
1,4-Dichlorobenzene-d4
Concen: 50.000 ug/l
RT: 12.006 min Scan# 11
Delta R.T. -0.000 min
Lab File: VX048733.D
Acq: 05 Dec 2025 14:13

Instrument :
MSVOA_X
ClientSampleId :
MW-14B-46-120325



Tgt Ion:	152	Resp:	131055
Ion	Ratio	Lower	Upper
152	100		
115	58.4	42.1	126.4
150	156.9	0.0	347.8





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

Report of Analysis

Client: JACOBS Engineering Group, Inc.
Project: Former Schlumberger STC PTC Site D3868221
Client Sample ID: MW-14B-46-120325RE
Lab Sample ID: Q3757-05RE
Analytical Method: 8260D
Sample Wt/Vol: 5 mL

Level: LOW
Final Vol: 5000 uL

Date Collected: 12/03/25
Date Received: 12/03/25
SDG No.: Q3757
Matrix: Water
% Solid: 0
Test: VOCMS Group3

CAS Number	Parameter	Conc.	Qua.	DF	MDL	LOQ / CRQL	Units	Date Ana.	BatchID
TARGETS									
75-01-4	Vinyl Chloride	0.26	U	1	0.26	1.00	ug/L	12/08/25 12:33	VX120825
75-35-4	1,1-Dichloroethene	0.23	U	1	0.23	1.00	ug/L	12/08/25 12:33	VX120825
75-34-3	1,1-Dichloroethane	0.23	U	1	0.23	1.00	ug/L	12/08/25 12:33	VX120825
156-59-2	cis-1,2-Dichloroethene	0.19	U	1	0.19	1.00	ug/L	12/08/25 12:33	VX120825
71-55-6	1,1,1-Trichloroethane	0.20	U	1	0.20	1.00	ug/L	12/08/25 12:33	VX120825
71-43-2	Benzene	0.15	U	1	0.15	1.00	ug/L	12/08/25 12:33	VX120825
107-06-2	1,2-Dichloroethane	0.22	U	1	0.22	1.00	ug/L	12/08/25 12:33	VX120825
79-01-6	Trichloroethene	0.090	U	1	0.090	1.00	ug/L	12/08/25 12:33	VX120825
79-00-5	1,1,2-Trichloroethane	0.21	U	1	0.21	1.00	ug/L	12/08/25 12:33	VX120825
127-18-4	Tetrachloroethene	0.23	U	1	0.23	1.00	ug/L	12/08/25 12:33	VX120825
SURROGATES									
17060-07-0	1,2-Dichloroethane-d4	57.7			70 (74) - 130 (125)	115%	SPK: 50		
1868-53-7	Dibromofluoromethane	38.4			70 (75) - 130 (124)	77%	SPK: 50		
2037-26-5	Toluene-d8	38.4			70 (86) - 130 (113)	77%	SPK: 50		
460-00-4	4-Bromofluorobenzene	56.9			70 (77) - 130 (121)	114%	SPK: 50		
INTERNAL STANDARDS									
		Area Count							
363-72-4	Pentafluorobenzene	125000							
540-36-3	1,4-Difluorobenzene	300000							
3114-55-4	Chlorobenzene-d5	268000							
3855-82-1	1,4-Dichlorobenzene-d4	140000							

U = Not Detected

LOQ = Limit of Quantitation

MDL = Method Detection Limit

LOD = Limit of Detection

E = Value Exceeds Calibration Range

Q = indicates LCS control criteria did not meet requirements

M = MS/MSD acceptance criteria did not meet requirements

J = Estimated Value

B = Analyte Found in Associated Method Blank

N = Presumptive Evidence of a Compound

* = Values outside of QC limits

D = Dilution

() = Laboratory InHouse Limit

A = Aldol-Condensation Reaction Products

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120825\
 Data File : VX048747.D
 Acq On : 08 Dec 2025 12:33
 Operator : JC/MD
 Sample : Q3757-05RE
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-14B-46-120325RE

Quant Time: Dec 09 04:08:38 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X120425W.M
 Quant Title : SW846 8260
 QLast Update : Fri Dec 05 03:27:34 2025
 Response via : Initial Calibration

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

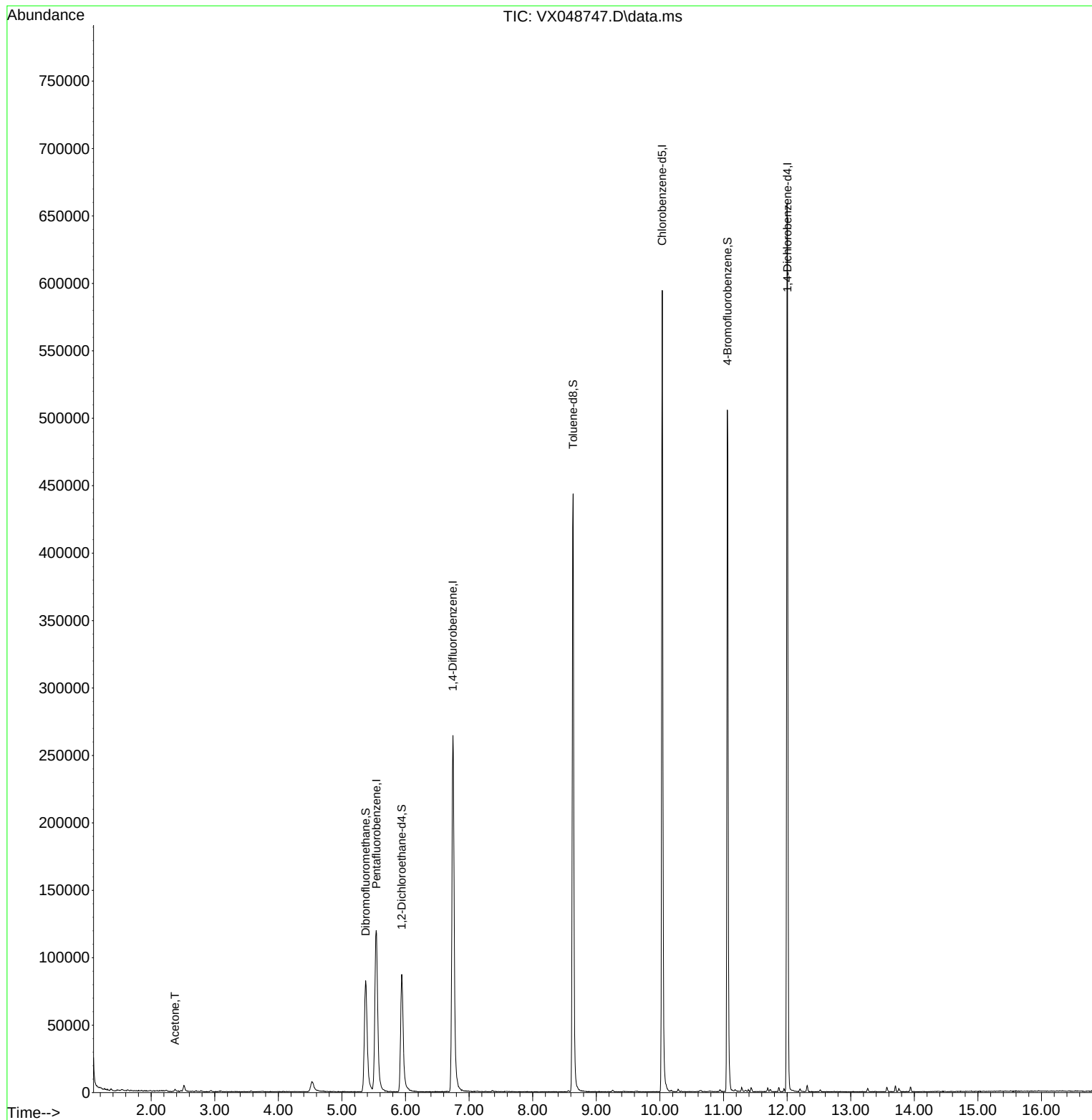
Internal Standards						
1) Pentafluorobenzene	5.544	168	124903	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.745	114	299763	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.037	117	268328	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.006	152	139875	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.940	65	96803	57.744	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	115.480%
35) Dibromofluoromethane	5.373	113	79281	38.412	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	76.820%
50) Toluene-d8	8.635	98	278126	38.442	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	76.880%#
62) 4-Bromofluorobenzene	11.061	95	141911	56.882	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	113.760%
Target Compounds						
16) Acetone	2.373	43	2472	3.852	ug/l	Qvalue # 72

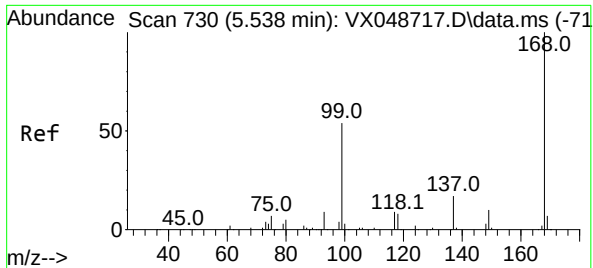
(#) = qualifier out of range (m) = manual integration (+) = signals summed

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120825\
Data File : VX048747.D
Acq On : 08 Dec 2025 12:33
Operator : JC/MD
Sample : Q3757-05RE
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-14B-46-120325RE

Quant Time: Dec 09 04:08:38 2025
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X120425W.M
Quant Title : SW846 8260
QLast Update : Fri Dec 05 03:27:34 2025
Response via : Initial Calibration

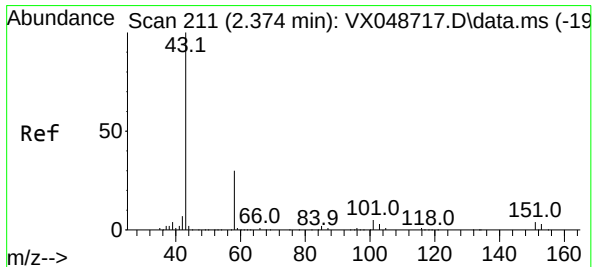
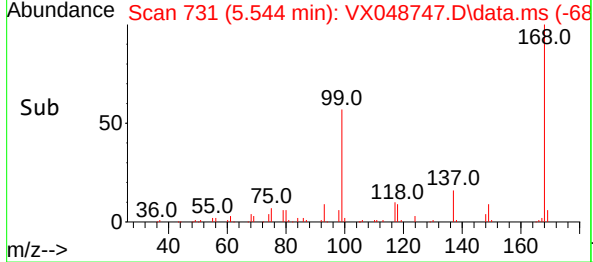
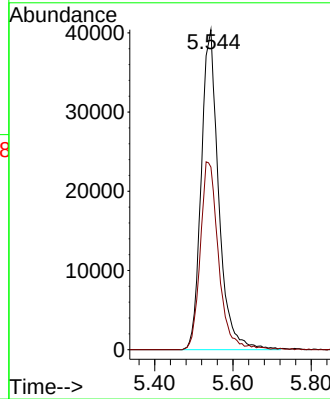
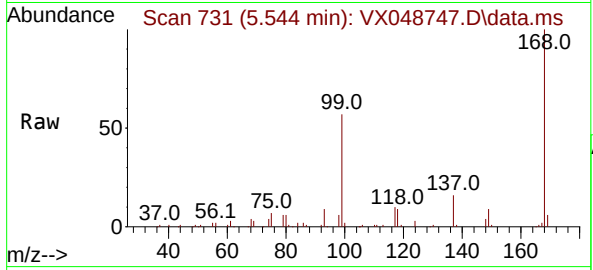




#1
Pentafluorobenzene
Concen: 50.000 ug/l
RT: 5.544 min Scan# 731
Delta R.T. 0.006 min
Lab File: VX048747.D
Acq: 08 Dec 2025 12:33

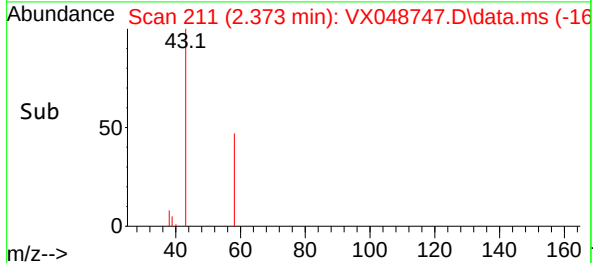
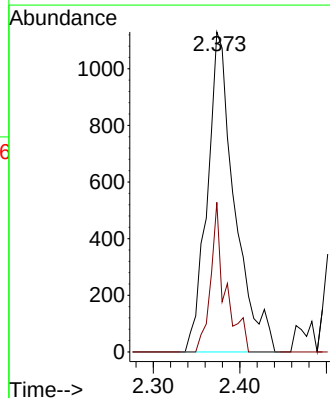
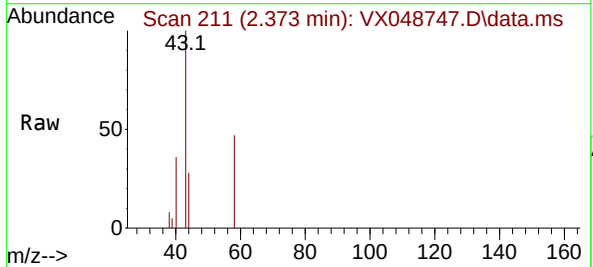
Instrument :
MSVOA_X
ClientSampleId :
MW-14B-46-120325RE

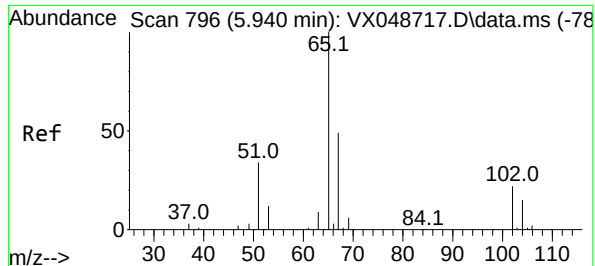
Tgt Ion:168 Resp: 124903
Ion Ratio Lower Upper
168 100
99 57.1 44.2 66.4



#16
Acetone
Concen: 3.852 ug/l
RT: 2.373 min Scan# 211
Delta R.T. -0.000 min
Lab File: VX048747.D
Acq: 08 Dec 2025 12:33

Tgt Ion: 43 Resp: 2472
Ion Ratio Lower Upper
43 100
58 46.7 25.0 37.4#





#33

1,2-Dichloroethane-d4

Concen: 57.744 ug/l

RT: 5.940 min Scan# 796

Delta R.T. -0.000 min

Lab File: VX048747.D

Acq: 08 Dec 2025 12:33

Instrument :

MSVOA_X

ClientSampleId :

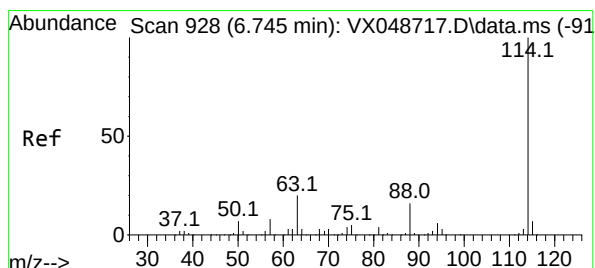
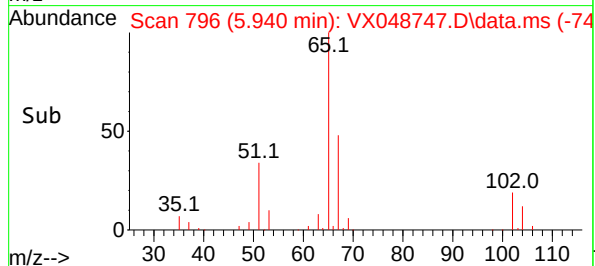
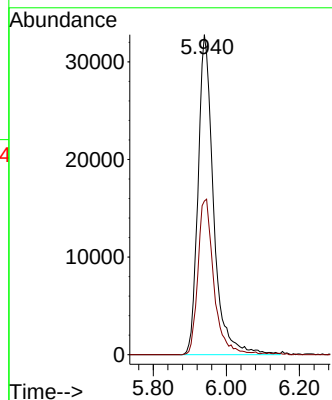
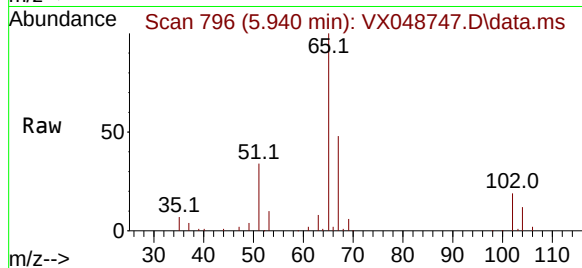
MW-14B-46-120325RE

Tgt Ion: 65 Resp: 96803

Ion Ratio Lower Upper

65 100

67 49.8 0.0 107.4



#34

1,4-Difluorobenzene

Concen: 50.000 ug/l

RT: 6.745 min Scan# 928

Delta R.T. -0.000 min

Lab File: VX048747.D

Acq: 08 Dec 2025 12:33

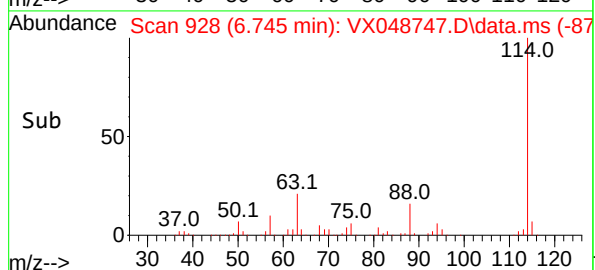
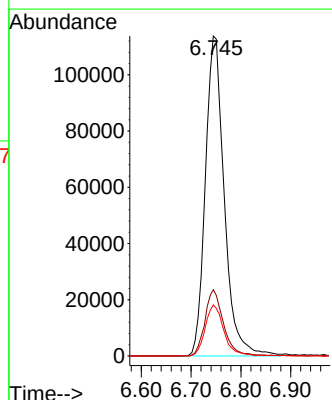
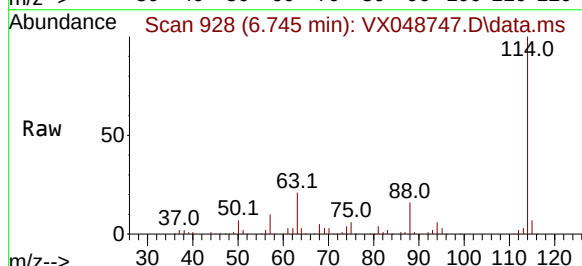
Tgt Ion: 114 Resp: 299763

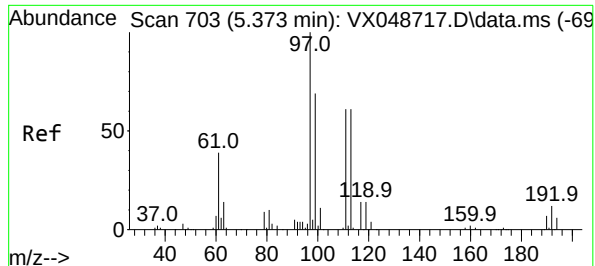
Ion Ratio Lower Upper

114 100

63 20.7 0.0 39.0

88 16.0 0.0 31.8





#35

Dibromofluoromethane

Concen: 38.412 ug/l

RT: 5.373 min Scan# 703

Delta R.T. -0.000 min

Lab File: VX048747.D

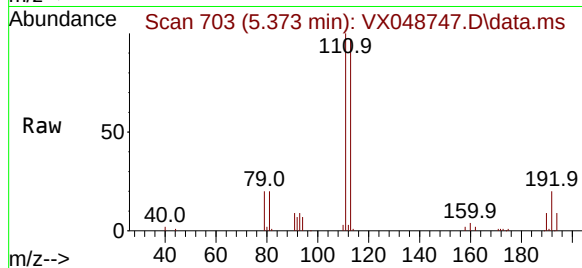
Acq: 08 Dec 2025 12:33

Instrument :

MSVOA_X

ClientSampleId :

MW-14B-46-120325RE



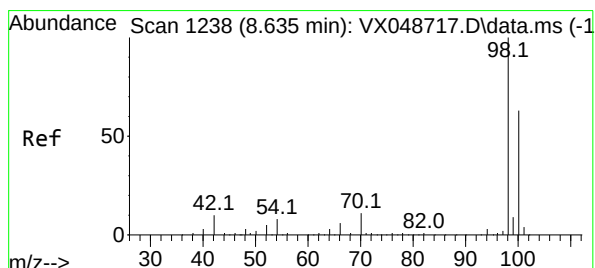
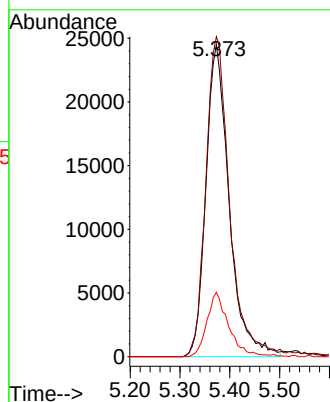
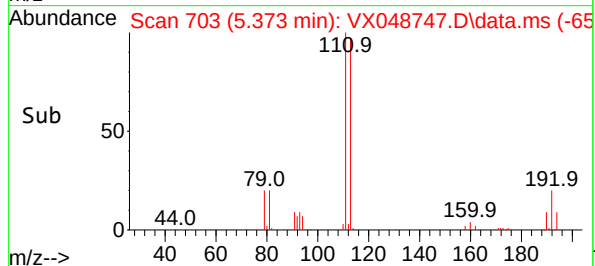
Tgt Ion:113 Resp: 79281

Ion Ratio Lower Upper

113 100

111 104.9 79.5 119.3

192 19.5 16.1 24.1



#50

Toluene-d8

Concen: 38.442 ug/l

RT: 8.635 min Scan# 1238

Delta R.T. -0.000 min

Lab File: VX048747.D

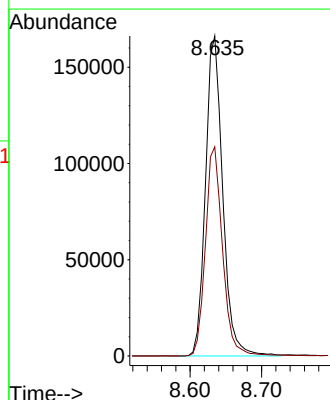
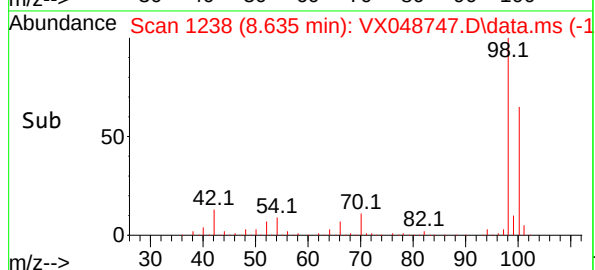
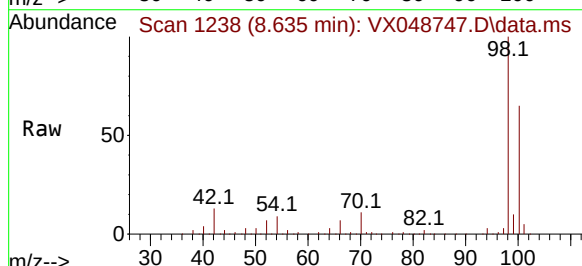
Acq: 08 Dec 2025 12:33

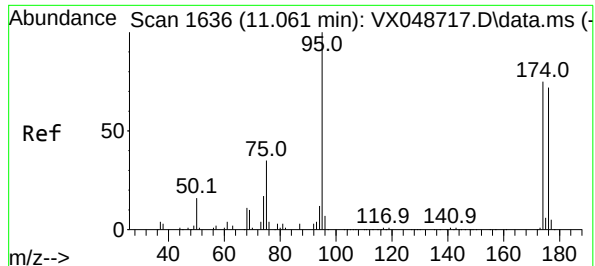
Tgt Ion: 98 Resp: 278126

Ion Ratio Lower Upper

98 100

100 65.2 53.4 80.0





#62

4-Bromofluorobenzene

Concen: 56.882 ug/l

RT: 11.061 min Scan# 1

Delta R.T. -0.000 min

Lab File: VX048747.D

Acq: 08 Dec 2025 12:33

Instrument :

MSVOA_X

ClientSampleId :

MW-14B-46-120325RE

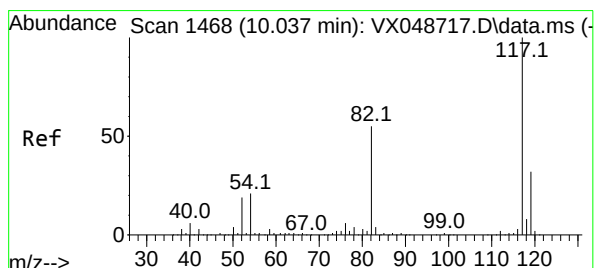
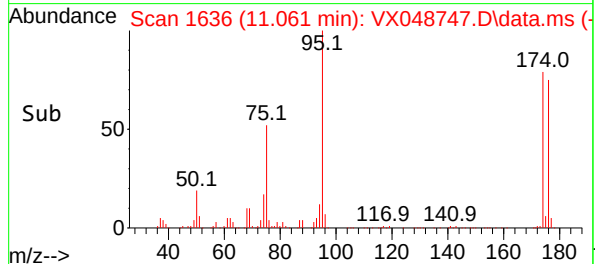
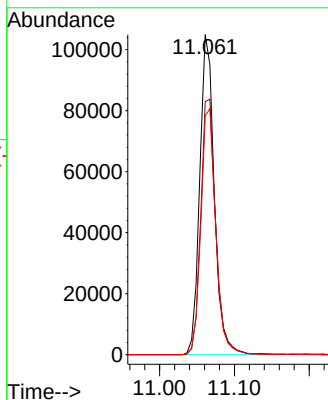
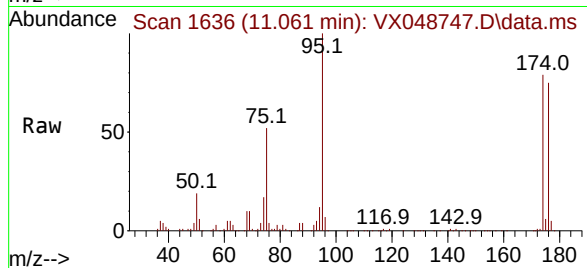
Tgt Ion: 95 Resp: 141911

Ion Ratio Lower Upper

95 100

174 82.4 0.0 157.8

176 79.0 0.0 154.0



#63

Chlorobenzene-d5

Concen: 50.000 ug/l

RT: 10.037 min Scan# 1468

Delta R.T. -0.000 min

Lab File: VX048747.D

Acq: 08 Dec 2025 12:33

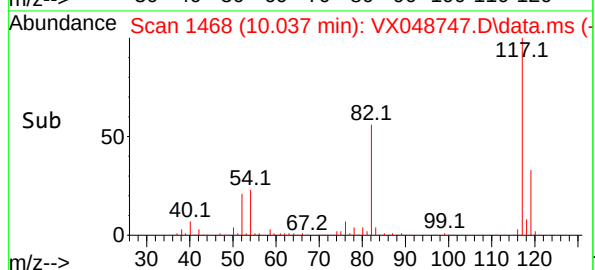
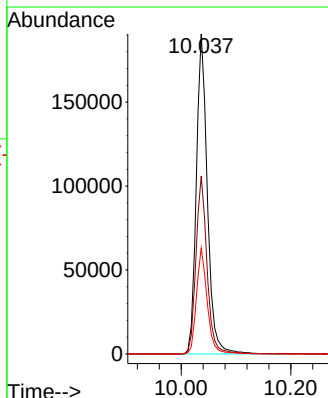
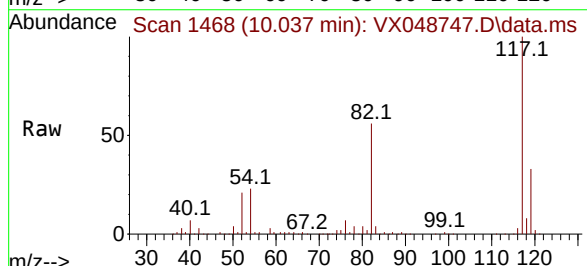
Tgt Ion: 117 Resp: 268328

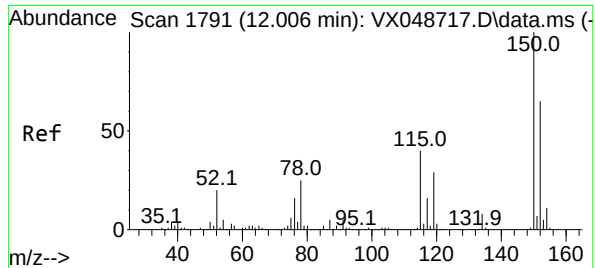
Ion Ratio Lower Upper

117 100

82 55.6 44.1 66.1

119 33.2 25.2 37.8





#72

1,4-Dichlorobenzene-d4

Concen: 50.000 ug/l

RT: 12.006 min Scan# 11

Delta R.T. -0.000 min

Lab File: VX048747.D

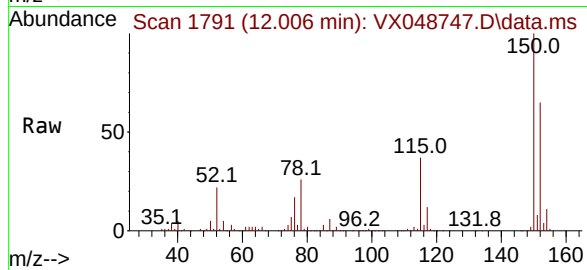
Acq: 08 Dec 2025 12:33

Instrument :

MSVOA_X

ClientSampleId :

MW-14B-46-120325RE



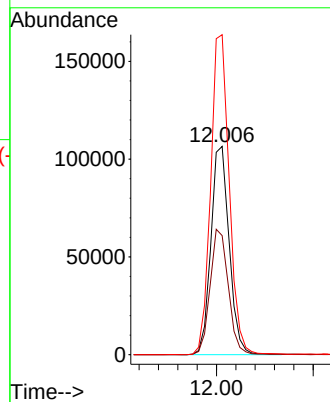
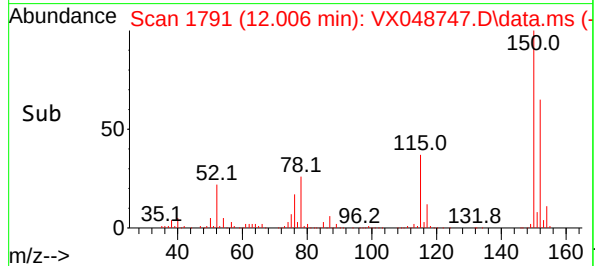
Tgt Ion:152 Resp: 139875

Ion Ratio Lower Upper

152 100

115 60.1 42.1 126.4

150 156.1 0.0 347.8





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

Report of Analysis

Client: JACOBS Engineering Group, Inc.
Project: Former Schlumberger STC PTC Site D3868221
Client Sample ID: RMW-05B-90.5-120325
Lab Sample ID: Q3757-07
Analytical Method: 8260D
Sample Wt/Vol: 5 mL

Level: LOW
Final Vol: 5000 uL

Date Collected: 12/03/25
Date Received: 12/03/25
SDG No.: Q3757
Matrix: Water
% Solid: 0
Test: VOCMS Group3

CAS Number	Parameter	Conc.	Qua.	DF	MDL	LOQ / CRQL	Units	Date Ana.	BatchID
TARGETS									
75-01-4	Vinyl Chloride	0.26	U	1	0.26	1.00	ug/L	12/05/25 14:33	VX120525
75-35-4	1,1-Dichloroethene	0.23	U	1	0.23	1.00	ug/L	12/05/25 14:33	VX120525
75-34-3	1,1-Dichloroethane	1.50		1	0.23	1.00	ug/L	12/05/25 14:33	VX120525
156-59-2	cis-1,2-Dichloroethene	13.1		1	0.19	1.00	ug/L	12/05/25 14:33	VX120525
71-55-6	1,1,1-Trichloroethane	0.20	U	1	0.20	1.00	ug/L	12/05/25 14:33	VX120525
71-43-2	Benzene	0.15	U	1	0.15	1.00	ug/L	12/05/25 14:33	VX120525
107-06-2	1,2-Dichloroethane	0.22	U	1	0.22	1.00	ug/L	12/05/25 14:33	VX120525
79-01-6	Trichloroethene	0.65	J	1	0.090	1.00	ug/L	12/05/25 14:33	VX120525
79-00-5	1,1,2-Trichloroethane	0.21	U	1	0.21	1.00	ug/L	12/05/25 14:33	VX120525
127-18-4	Tetrachloroethene	0.23	U	1	0.23	1.00	ug/L	12/05/25 14:33	VX120525
SURROGATES									
17060-07-0	1,2-Dichloroethane-d4	51.3			70 (74) - 130 (125)	103%	SPK: 50		
1868-53-7	Dibromofluoromethane	47.1			70 (75) - 130 (124)	94%	SPK: 50		
2037-26-5	Toluene-d8	43.9			70 (86) - 130 (113)	88%	SPK: 50		
460-00-4	4-Bromofluorobenzene	45.8			70 (77) - 130 (121)	92%	SPK: 50		
INTERNAL STANDARDS									
		Area Count							
363-72-4	Pentafluorobenzene	132000							
540-36-3	1,4-Difluorobenzene	223000							
3114-55-4	Chlorobenzene-d5	207000							
3855-82-1	1,4-Dichlorobenzene-d4	103000							

U = Not Detected

LOQ = Limit of Quantitation

MDL = Method Detection Limit

LOD = Limit of Detection

E = Value Exceeds Calibration Range

Q = indicates LCS control criteria did not meet requirements

M = MS/MSD acceptance criteria did not meet requirements

J = Estimated Value

B = Analyte Found in Associated Method Blank

N = Presumptive Evidence of a Compound

* = Values outside of QC limits

D = Dilution

() = Laboratory InHouse Limit

A = Aldol-Condensation Reaction Products

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120525\
 Data File : VX048734.D
 Acq On : 05 Dec 2025 14:33
 Operator : JC/MD
 Sample : Q3757-07
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 RMW-05B-90.5-120325

Quant Time: Dec 05 22:05:03 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X120425W.M
 Quant Title : SW846 8260
 QLast Update : Fri Dec 05 03:27:34 2025
 Response via : Initial Calibration

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

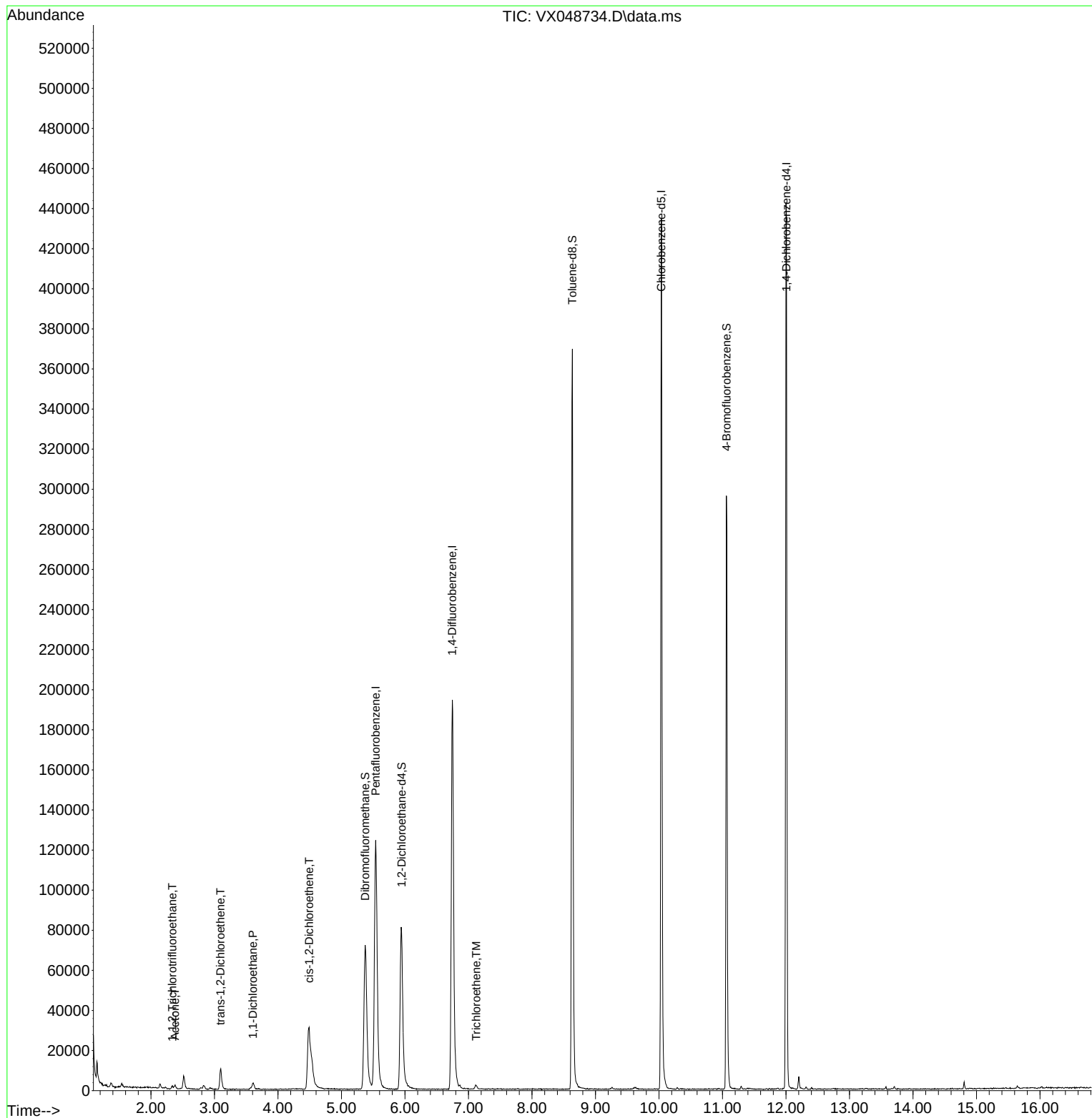
Internal Standards						
1) Pentafluorobenzene	5.538	168	132392	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.745	114	222932	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.037	117	207009	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.006	152	102583	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.946	65	91135	51.288	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	= 102.580%	
35) Dibromofluoromethane	5.373	113	72316	47.112	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 94.220%	
50) Toluene-d8	8.635	98	236336	43.924	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	= 87.840%#	
62) 4-Bromofluorobenzene	11.061	95	84945	45.783	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 91.560%	
Target Compounds						Qvalue
9) 1,1,2-Trichlorotrifluo...	2.337	101	558	0.370	ug/l #	83
16) Acetone	2.380	43	3009	4.423	ug/l #	87
21) trans-1,2-Dichloroethene	3.093	96	5495	3.502	ug/l #	79
24) 1,1-Dichloroethane	3.605	63	3969	1.475	ug/l	97
27) cis-1,2-Dichloroethene	4.489	96	23281	13.149	ug/l	90
44) Trichloroethene	7.123	130	1059	0.651	ug/l	91

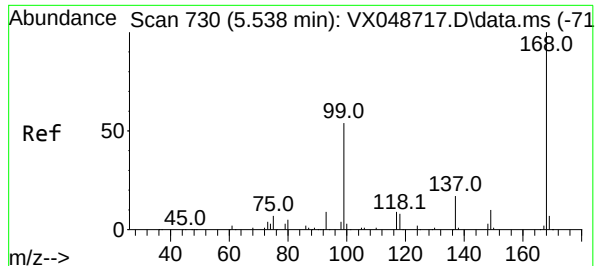
(#) = qualifier out of range (m) = manual integration (+) = signals summed

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120525\
Data File : VX048734.D
Acq On : 05 Dec 2025 14:33
Operator : JC/MD
Sample : Q3757-07
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
RMW-05B-90.5-120325

Quant Time: Dec 05 22:05:03 2025
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X120425W.M
Quant Title : SW846 8260
QLast Update : Fri Dec 05 03:27:34 2025
Response via : Initial Calibration

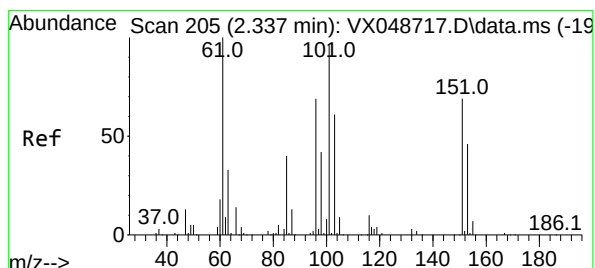
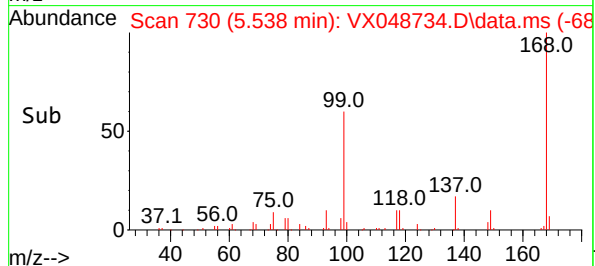
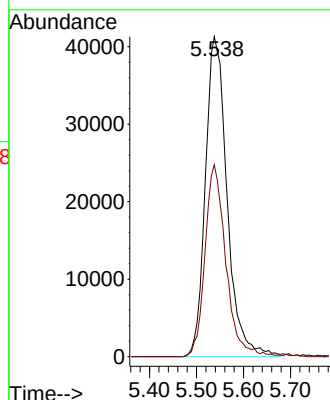
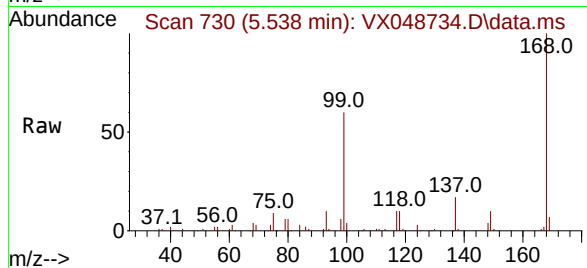




#1
Pentafluorobenzene
Concen: 50.000 ug/l
RT: 5.538 min Scan# 71
Delta R.T. -0.000 min
Lab File: VX048734.D
Acq: 05 Dec 2025 14:33

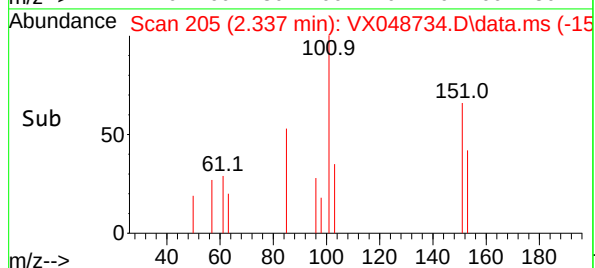
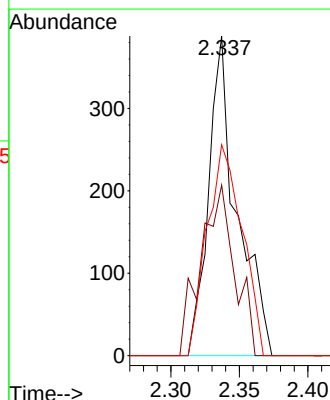
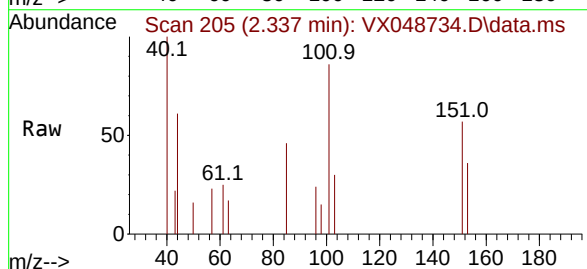
Instrument :
MSVOA_X
ClientSampleId :
RMW-05B-90.5-120325

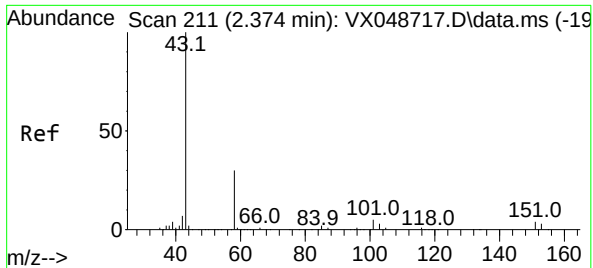
Tgt Ion:168 Resp: 132392
Ion Ratio Lower Upper
168 100
99 59.9 44.2 66.4



#9
1,1,2-Trichlorotrifluoroethane
Concen: 0.370 ug/l
RT: 2.337 min Scan# 205
Delta R.T. 0.000 min
Lab File: VX048734.D
Acq: 05 Dec 2025 14:33

Tgt Ion:101 Resp: 558
Ion Ratio Lower Upper
101 100
85 64.0 34.2 51.2#
151 81.5 60.2 90.4

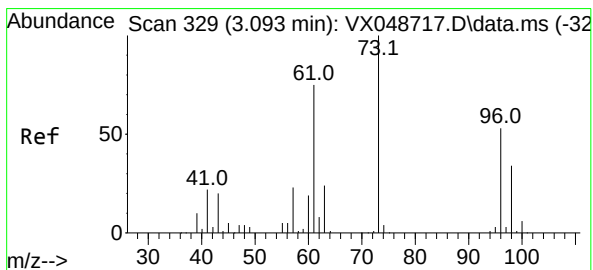
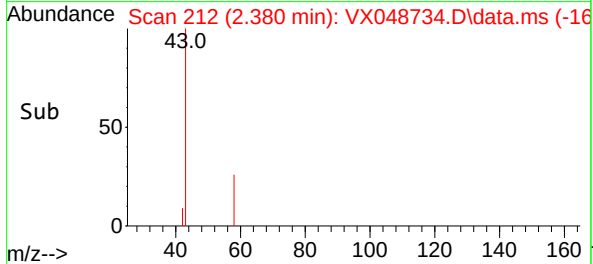
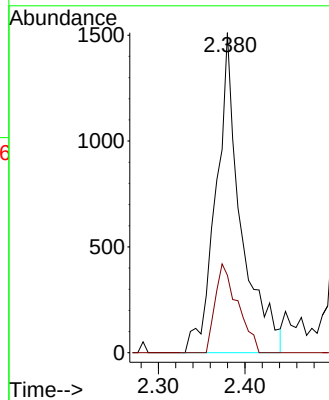
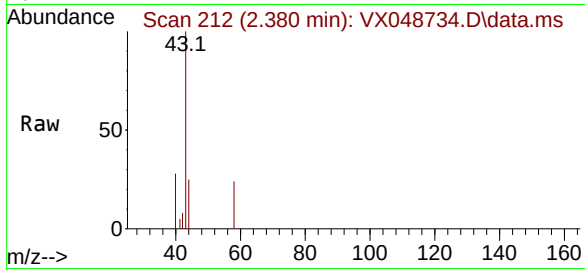




#16
Acetone
Concen: 4.423 ug/l
RT: 2.380 min Scan# 211
Delta R.T. 0.006 min
Lab File: VX048734.D
Acq: 05 Dec 2025 14:33

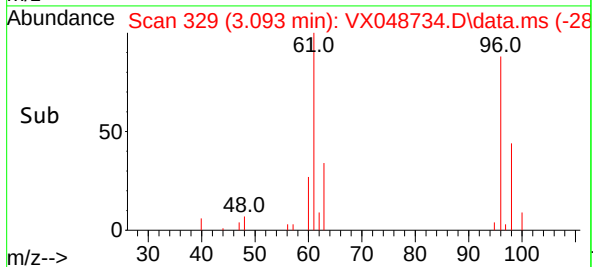
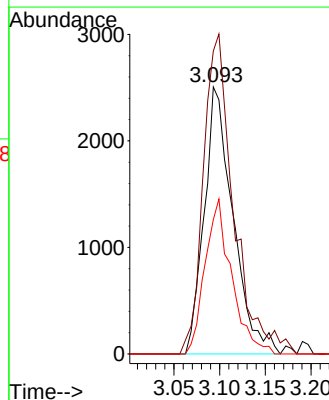
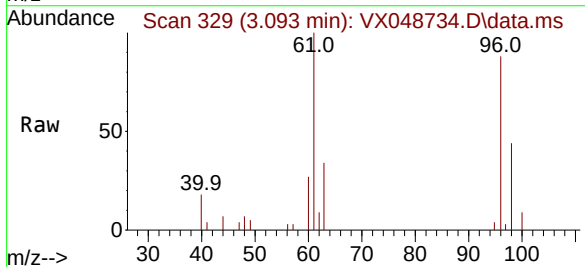
Instrument :
MSVOA_X
ClientSampleId :
RMW-05B-90.5-120325

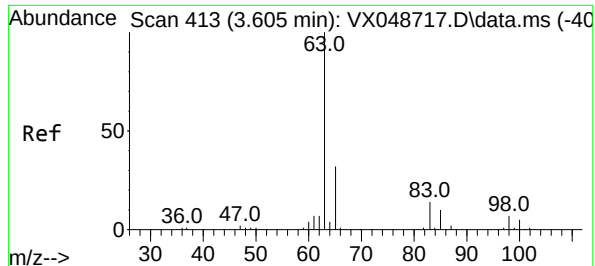
Tgt Ion: 43 Resp: 3009
Ion Ratio Lower Upper
43 100
58 24.2 25.0 37.4#



#21
trans-1,2-Dichloroethene
Concen: 3.502 ug/l
RT: 3.093 min Scan# 329
Delta R.T. 0.000 min
Lab File: VX048734.D
Acq: 05 Dec 2025 14:33

Tgt Ion: 96 Resp: 5495
Ion Ratio Lower Upper
96 100
61 113.4 112.8 169.2
98 50.3 51.0 76.4#

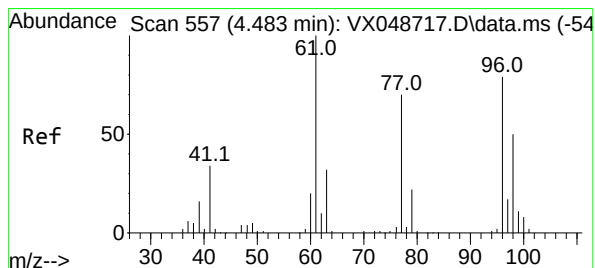
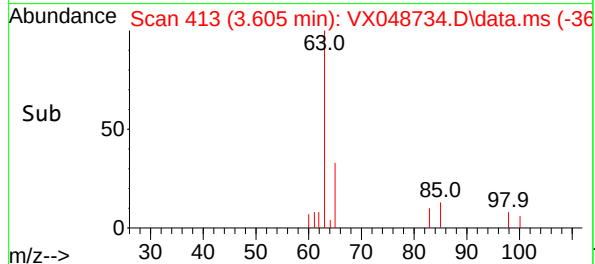
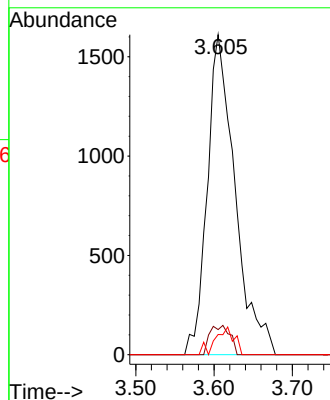
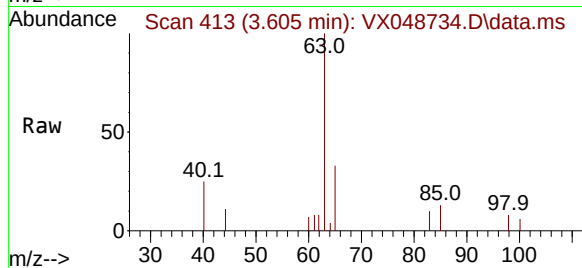




#24
 1,1-Dichloroethane
 Concen: 1.475 ug/l
 RT: 3.605 min Scan# 413
 Delta R.T. 0.000 min
 Lab File: VX048734.D
 Acq: 05 Dec 2025 14:33

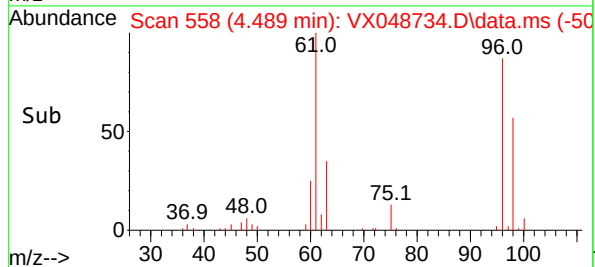
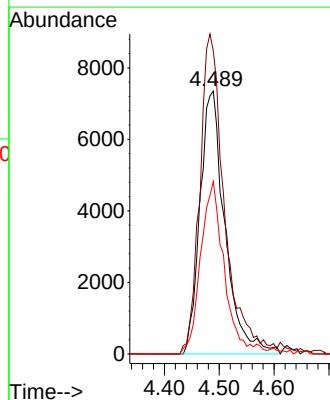
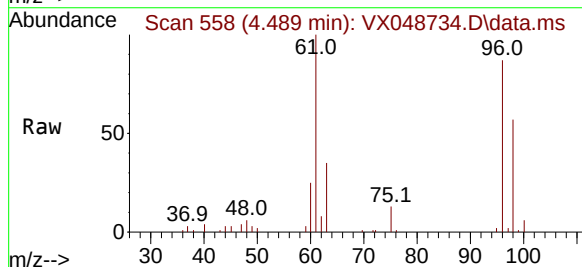
Instrument :
 MSVOA_X
 ClientSampleId :
 RMW-05B-90.5-120325

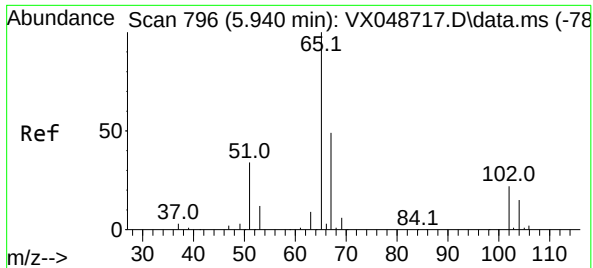
Tgt Ion:	63	Resp:	3969
Ion	Ratio	Lower	Upper
63	100		
98	7.9	3.5	10.6
100	6.3	2.5	7.4



#27
 cis-1,2-Dichloroethene
 Concen: 13.149 ug/l
 RT: 4.489 min Scan# 558
 Delta R.T. 0.006 min
 Lab File: VX048734.D
 Acq: 05 Dec 2025 14:33

Tgt Ion:	96	Resp:	23281
Ion	Ratio	Lower	Upper
96	100		
61	119.9	0.0	271.0
98	61.8	0.0	130.2





#33

1,2-Dichloroethane-d4

Concen: 51.288 ug/l

RT: 5.946 min Scan# 796

Delta R.T. 0.006 min

Lab File: VX048734.D

Acq: 05 Dec 2025 14:33

Instrument :

MSVOA_X

ClientSampleId :

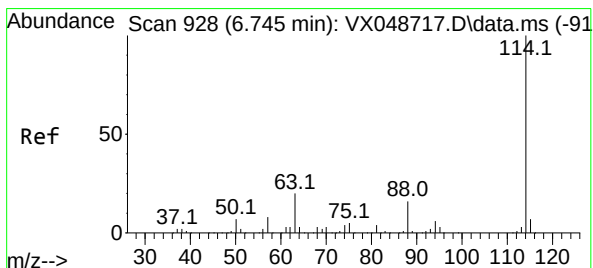
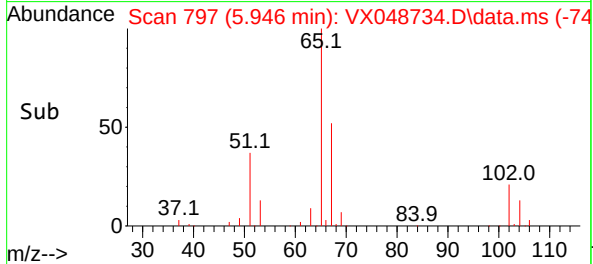
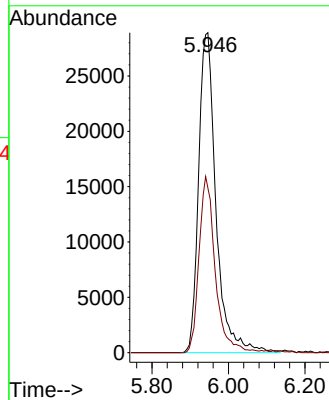
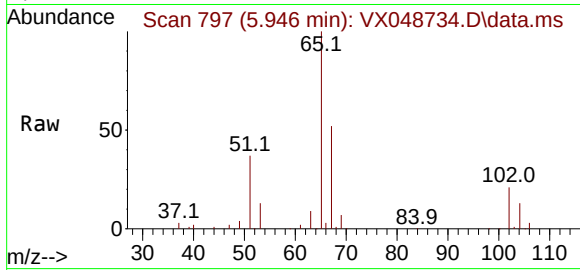
RMW-05B-90.5-120325

Tgt Ion: 65 Resp: 91135

Ion Ratio Lower Upper

65 100

67 51.4 0.0 107.4



#34

1,4-Difluorobenzene

Concen: 50.000 ug/l

RT: 6.745 min Scan# 928

Delta R.T. 0.000 min

Lab File: VX048734.D

Acq: 05 Dec 2025 14:33

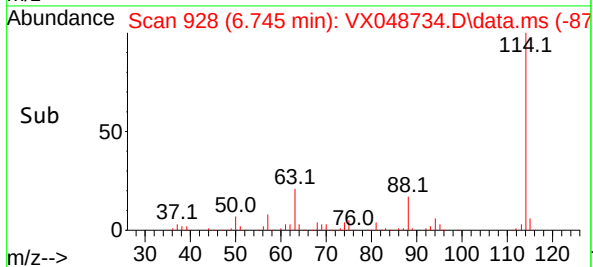
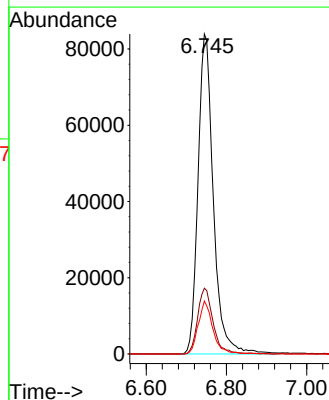
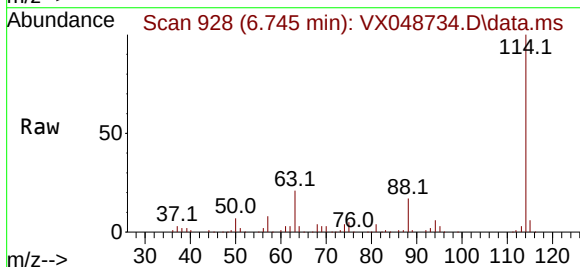
Tgt Ion: 114 Resp: 222932

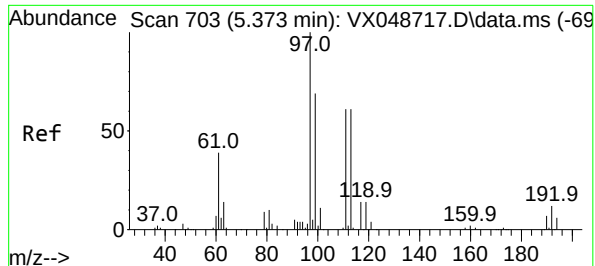
Ion Ratio Lower Upper

114 100

63 20.6 0.0 39.0

88 16.6 0.0 31.8





#35

Dibromofluoromethane

Concen: 47.112 ug/l

RT: 5.373 min Scan# 703

Delta R.T. 0.000 min

Lab File: VX048734.D

Acq: 05 Dec 2025 14:33

Instrument :

MSVOA_X

ClientSampleId :

RMW-05B-90.5-120325

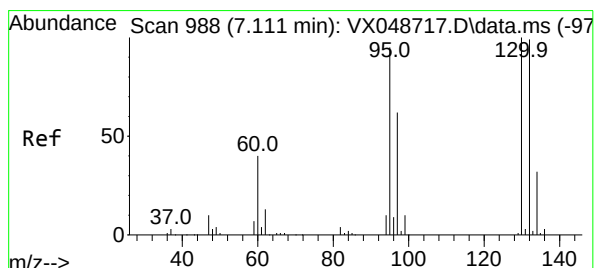
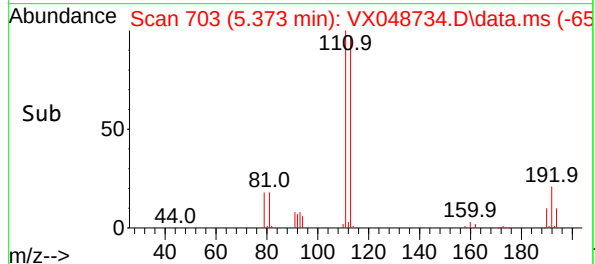
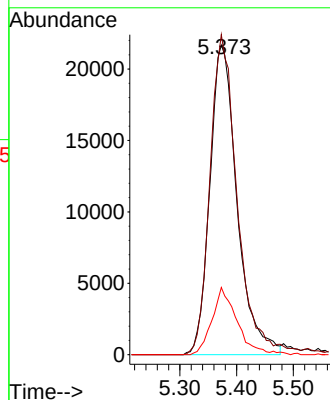
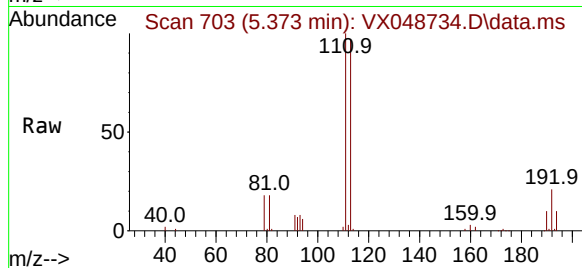
Tgt Ion:113 Resp: 72316

Ion Ratio Lower Upper

113 100

111 103.5 79.5 119.3

192 19.6 16.1 24.1



#44

Trichloroethene

Concen: 0.651 ug/l

RT: 7.123 min Scan# 990

Delta R.T. 0.012 min

Lab File: VX048734.D

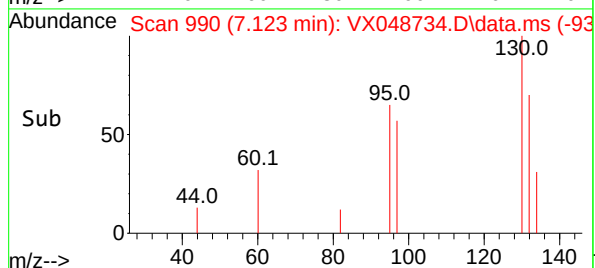
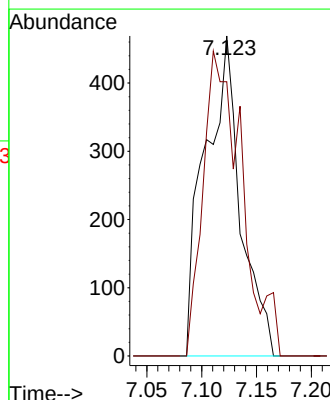
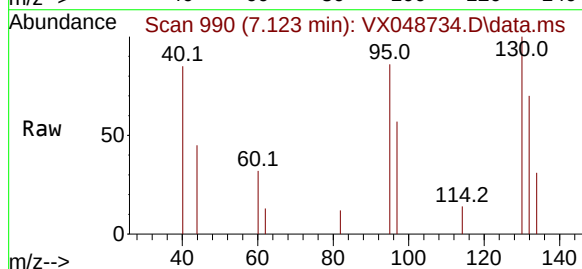
Acq: 05 Dec 2025 14:33

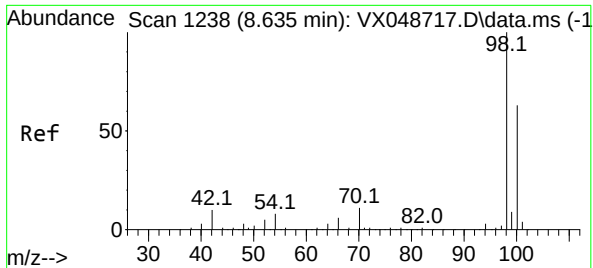
Tgt Ion:130 Resp: 1059

Ion Ratio Lower Upper

130 100

95 85.7 0.0 189.6

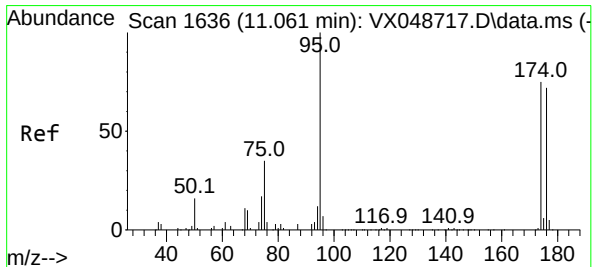
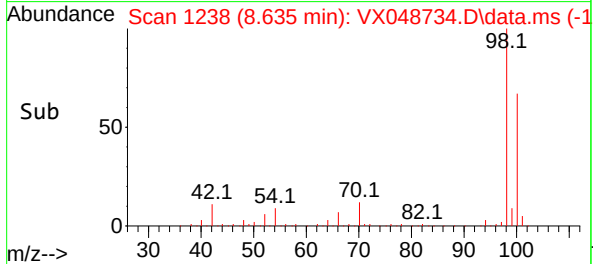
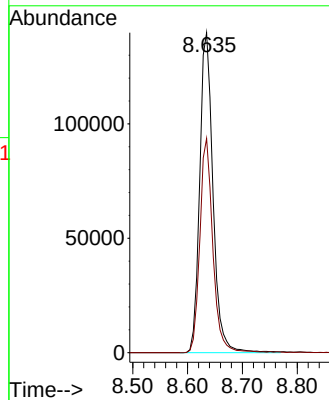
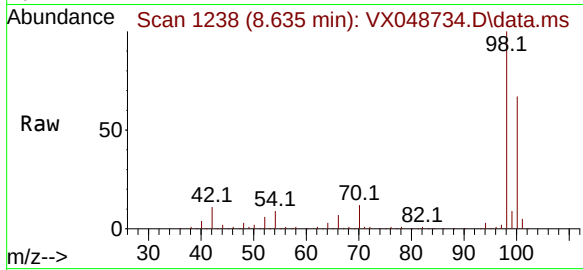




#50
Toluene-d8
Concen: 43.924 ug/l
RT: 8.635 min Scan# 11
Delta R.T. 0.000 min
Lab File: VX048734.D
Acq: 05 Dec 2025 14:33

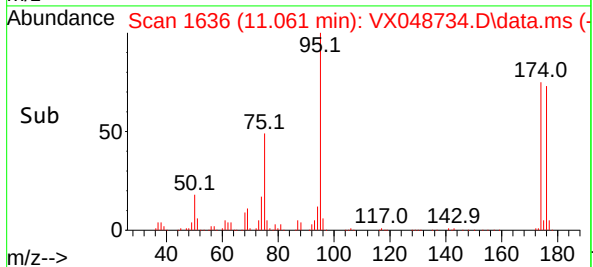
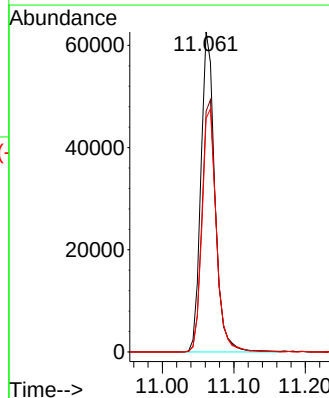
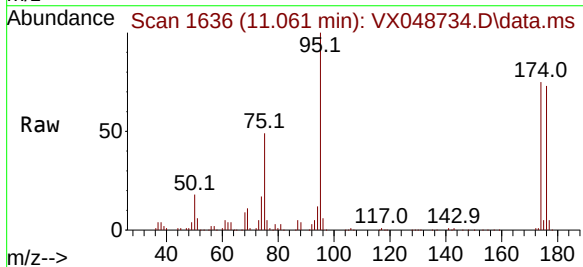
Instrument :
MSVOA_X
ClientSampleId :
RMW-05B-90.5-120325

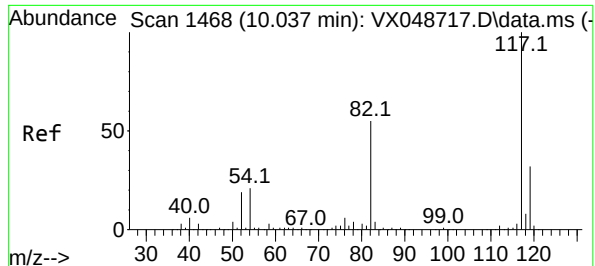
Tgt Ion: 98 Resp: 236336
Ion Ratio Lower Upper
98 100
100 65.4 53.4 80.0



#62
4-Bromofluorobenzene
Concen: 45.783 ug/l
RT: 11.061 min Scan# 1636
Delta R.T. 0.000 min
Lab File: VX048734.D
Acq: 05 Dec 2025 14:33

Tgt Ion: 95 Resp: 84945
Ion Ratio Lower Upper
95 100
174 80.1 0.0 157.8
176 78.0 0.0 154.0





#63

Chlorobenzene-d5

Concen: 50.000 ug/l

RT: 10.037 min Scan# 1468

Delta R.T. 0.000 min

Lab File: VX048734.D

Acq: 05 Dec 2025 14:33

Instrument :

MSVOA_X

ClientSampleId :

RMW-05B-90.5-120325

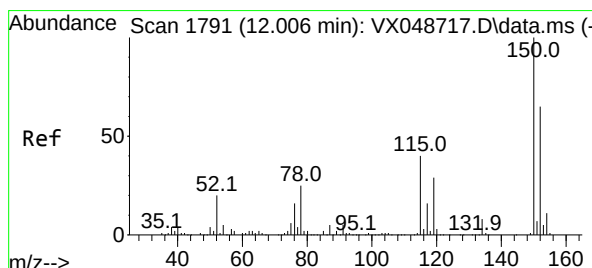
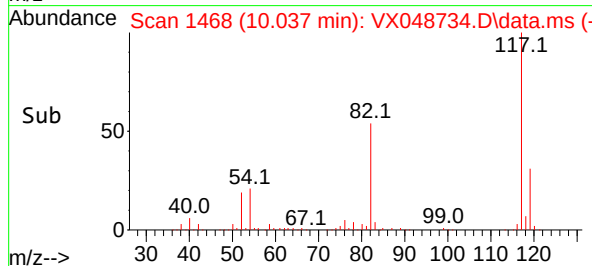
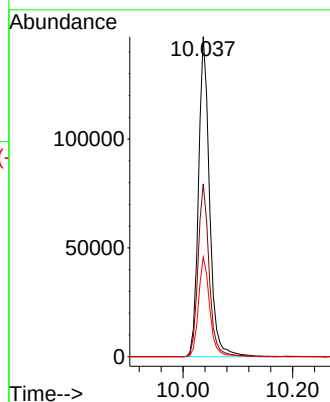
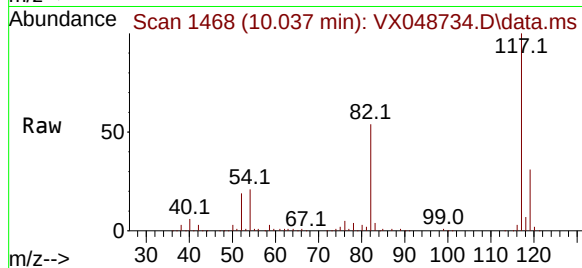
Tgt Ion:117 Resp: 207009

Ion Ratio Lower Upper

117 100

82 54.0 44.1 66.1

119 31.1 25.2 37.8



#72

1,4-Dichlorobenzene-d4

Concen: 50.000 ug/l

RT: 12.006 min Scan# 1791

Delta R.T. 0.000 min

Lab File: VX048734.D

Acq: 05 Dec 2025 14:33

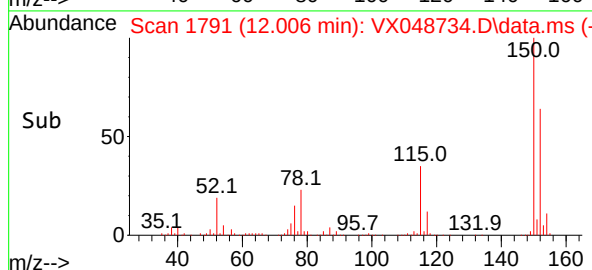
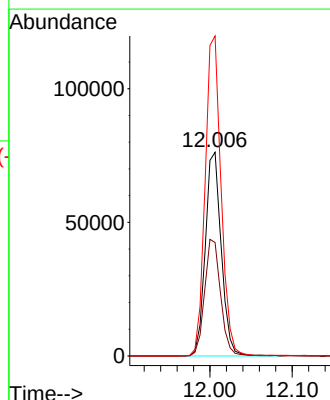
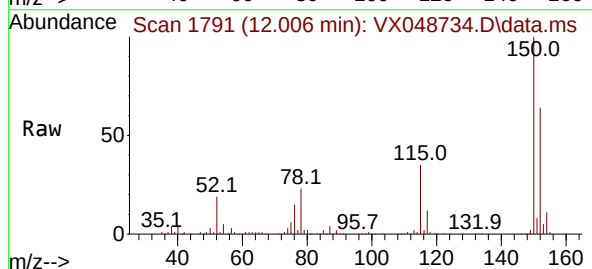
Tgt Ion:152 Resp: 102583

Ion Ratio Lower Upper

152 100

115 57.5 42.1 126.4

150 156.8 0.0 347.8





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

Report of Analysis

Client: JACOBS Engineering Group, Inc.
Project: Former Schlumberger STC PTC Site D3868221
Client Sample ID: OW-05B-72-120325
Lab Sample ID: Q3757-09
Analytical Method: 8260D
Sample Wt/Vol: 5 mL

Level: LOW
Final Vol: 5000 uL

Date Collected: 12/03/25
Date Received: 12/03/25
SDG No.: Q3757
Matrix: Water
% Solid: 0
Test: VOCMS Group3

CAS Number	Parameter	Conc.	Qua.	DF	MDL	LOQ / CRQL	Units	Date Ana.	BatchID
TARGETS									
75-01-4	Vinyl Chloride	0.26	U	1	0.26	1.00	ug/L	12/05/25 14:54	VX120525
75-35-4	1,1-Dichloroethene	0.85	J	1	0.23	1.00	ug/L	12/05/25 14:54	VX120525
75-34-3	1,1-Dichloroethane	0.23	U	1	0.23	1.00	ug/L	12/05/25 14:54	VX120525
156-59-2	cis-1,2-Dichloroethene	2.20		1	0.19	1.00	ug/L	12/05/25 14:54	VX120525
71-55-6	1,1,1-Trichloroethane	0.20	U	1	0.20	1.00	ug/L	12/05/25 14:54	VX120525
71-43-2	Benzene	0.15	U	1	0.15	1.00	ug/L	12/05/25 14:54	VX120525
107-06-2	1,2-Dichloroethane	0.22	U	1	0.22	1.00	ug/L	12/05/25 14:54	VX120525
79-01-6	Trichloroethene	0.44	J	1	0.090	1.00	ug/L	12/05/25 14:54	VX120525
79-00-5	1,1,2-Trichloroethane	0.21	U	1	0.21	1.00	ug/L	12/05/25 14:54	VX120525
127-18-4	Tetrachloroethene	0.23	U	1	0.23	1.00	ug/L	12/05/25 14:54	VX120525
SURROGATES									
17060-07-0	1,2-Dichloroethane-d4	58.8			70 (74) - 130 (125)	118%	SPK: 50		
1868-53-7	Dibromofluoromethane	39.7			70 (75) - 130 (124)	79%	SPK: 50		
2037-26-5	Toluene-d8	38.2			70 (86) - 130 (113)	76%	SPK: 50		
460-00-4	4-Bromofluorobenzene	45.0			70 (77) - 130 (121)	90%	SPK: 50		
INTERNAL STANDARDS									
		Area Count							
363-72-4	Pentafluorobenzene	129000							
540-36-3	1,4-Difluorobenzene	275000							
3114-55-4	Chlorobenzene-d5	209000							
3855-82-1	1,4-Dichlorobenzene-d4	103000							

U = Not Detected

LOQ = Limit of Quantitation

MDL = Method Detection Limit

LOD = Limit of Detection

E = Value Exceeds Calibration Range

Q = indicates LCS control criteria did not meet requirements

M = MS/MSD acceptance criteria did not meet requirements

J = Estimated Value

B = Analyte Found in Associated Method Blank

N = Presumptive Evidence of a Compound

* = Values outside of QC limits

D = Dilution

() = Laboratory InHouse Limit

A = Aldol-Condensation Reaction Products

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120525\
 Data File : VX048735.D
 Acq On : 05 Dec 2025 14:54
 Operator : JC/MD
 Sample : Q3757-09
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 OW-05B-72-120325

Quant Time: Dec 05 22:05:28 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X120425W.M
 Quant Title : SW846 8260
 QLast Update : Fri Dec 05 03:27:34 2025
 Response via : Initial Calibration

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

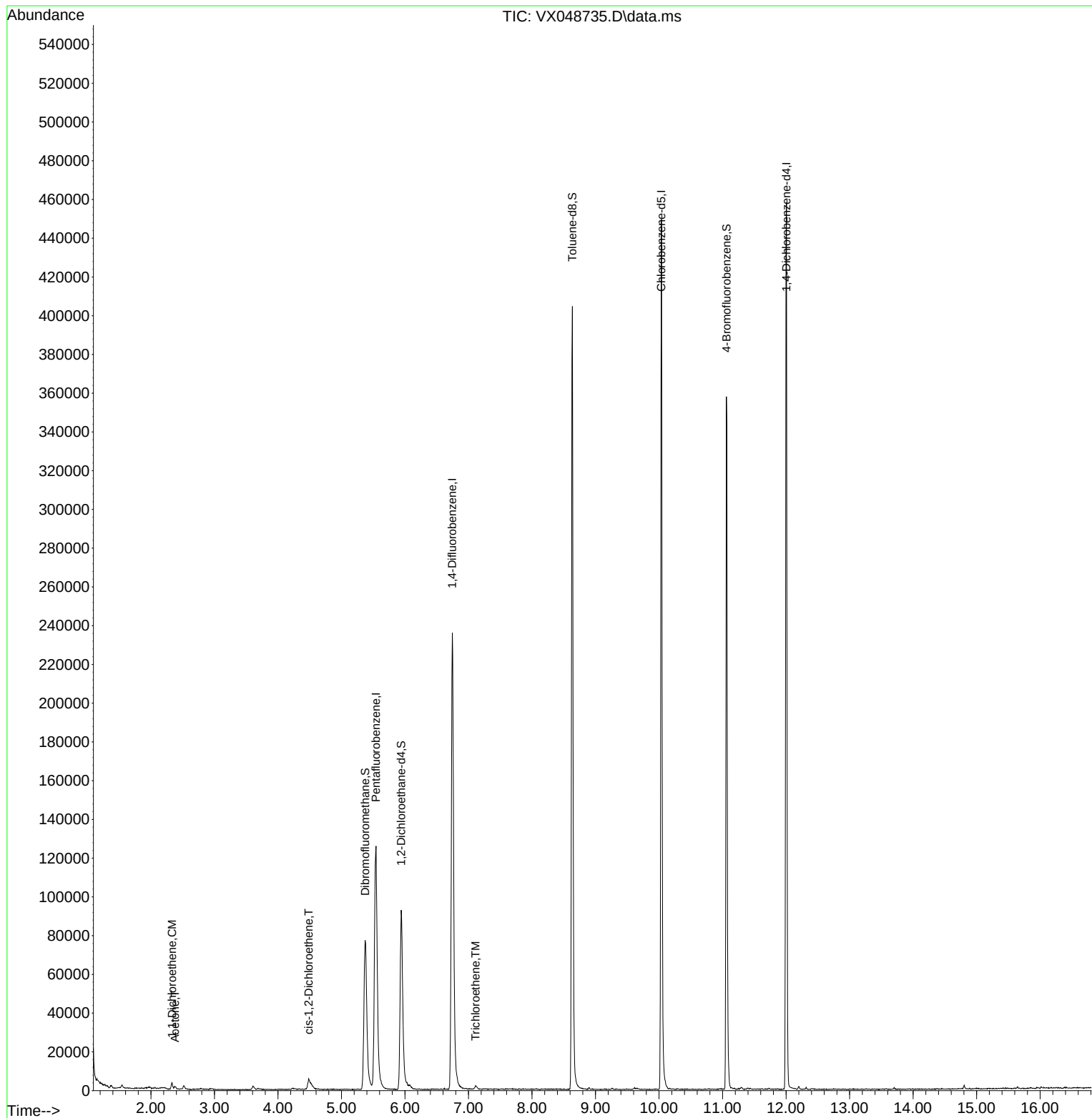
Internal Standards						
1) Pentafluorobenzene	5.544	168	129477	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.745	114	275147	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.037	117	208806	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.006	152	103213	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.940	65	102254	58.840	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	117.680%#
35) Dibromofluoromethane	5.373	113	75188	39.688	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	79.380%
50) Toluene-d8	8.635	98	253556	38.182	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	76.360%#
62) 4-Bromofluorobenzene	11.061	95	103159	45.048	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	90.100%
Target Compounds						
12) 1,1-Dichloroethene	2.331	96	1284	0.854	ug/l #	81
16) Acetone	2.380	43	2234	3.358	ug/l	91
27) cis-1,2-Dichloroethene	4.483	96	3749	2.165	ug/l	92
44) Trichloroethene	7.110	130	883	0.440	ug/l	86

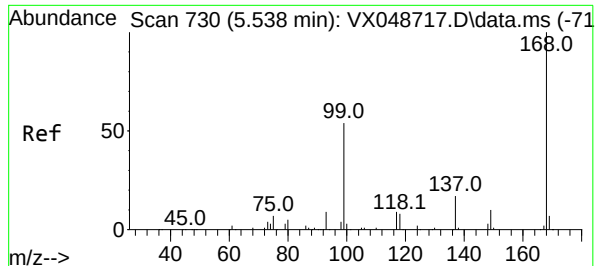
(#) = qualifier out of range (m) = manual integration (+) = signals summed

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120525\
Data File : VX048735.D
Acq On : 05 Dec 2025 14:54
Operator : JC/MD
Sample : Q3757-09
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
OW-05B-72-120325

Quant Time: Dec 05 22:05:28 2025
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X120425W.M
Quant Title : SW846 8260
QLast Update : Fri Dec 05 03:27:34 2025
Response via : Initial Calibration

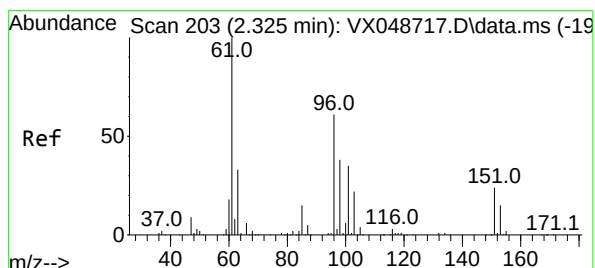
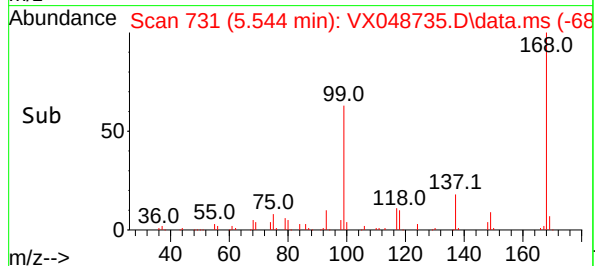
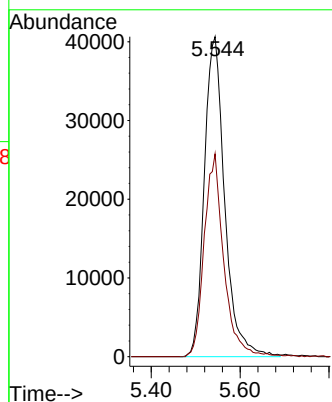
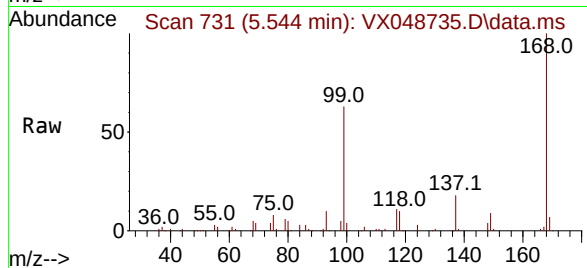




#1
Pentafluorobenzene
Concen: 50.000 ug/l
RT: 5.544 min Scan# 71
Delta R.T. 0.006 min
Lab File: VX048735.D
Acq: 05 Dec 2025 14:54

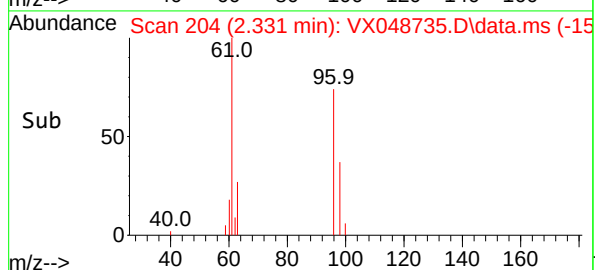
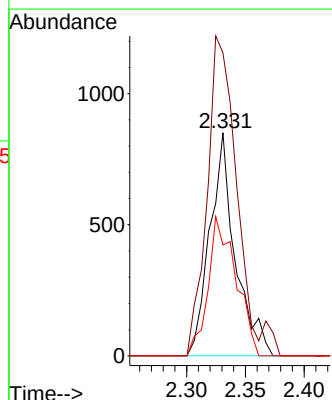
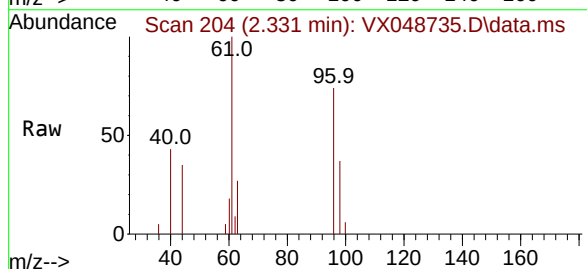
Instrument :
MSVOA_X
ClientSampleId :
OW-05B-72-120325

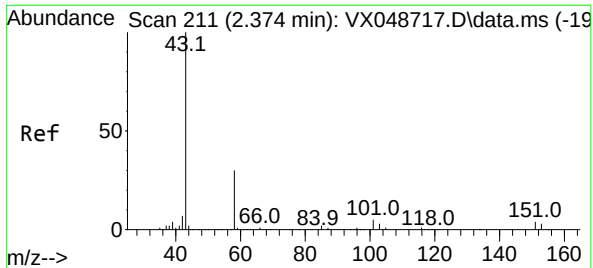
Tgt Ion:168 Resp: 129477
Ion Ratio Lower Upper
168 100
99 63.4 44.2 66.4



#12
1,1-Dichloroethene
Concen: 0.854 ug/l
RT: 2.331 min Scan# 204
Delta R.T. 0.006 min
Lab File: VX048735.D
Acq: 05 Dec 2025 14:54

Tgt Ion: 96 Resp: 1284
Ion Ratio Lower Upper
96 100
61 135.8 130.7 196.1
98 49.7 49.8 74.8#

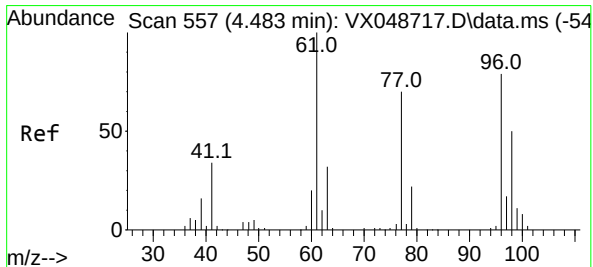
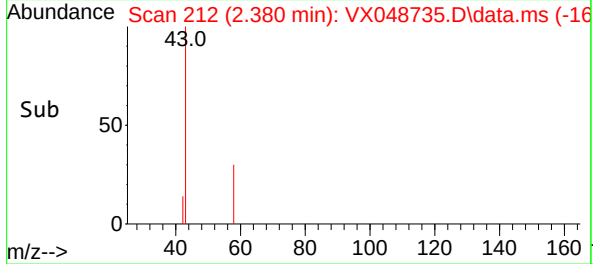
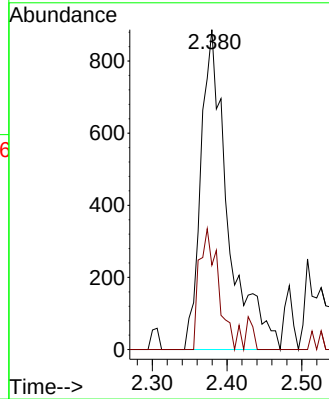
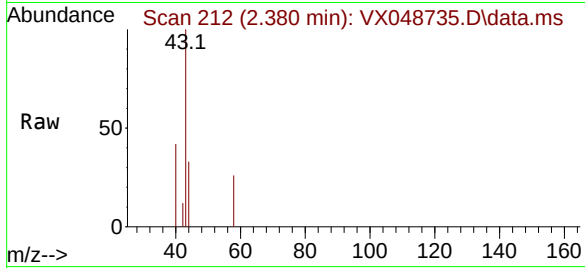




#16
Acetone
Concen: 3.358 ug/l
RT: 2.380 min Scan# 211
Delta R.T. 0.006 min
Lab File: VX048735.D
Acq: 05 Dec 2025 14:54

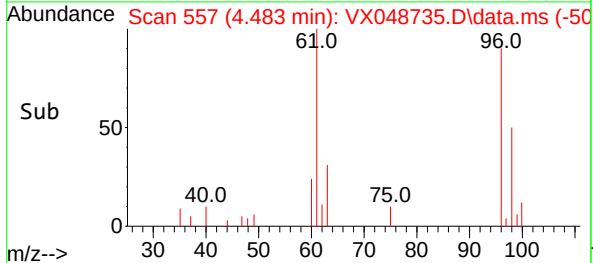
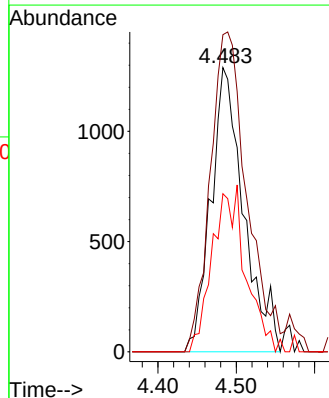
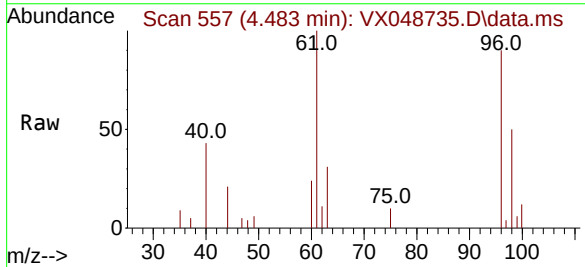
Instrument :
MSVOA_X
ClientSampleId :
OW-05B-72-120325

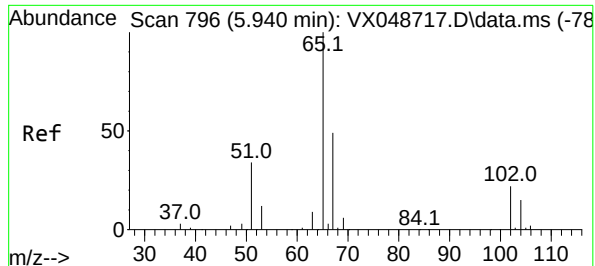
Tgt Ion: 43 Resp: 2234
Ion Ratio Lower Upper
43 100
58 26.4 25.0 37.4



#27
cis-1,2-Dichloroethene
Concen: 2.165 ug/l
RT: 4.483 min Scan# 557
Delta R.T. -0.000 min
Lab File: VX048735.D
Acq: 05 Dec 2025 14:54

Tgt Ion: 96 Resp: 3749
Ion Ratio Lower Upper
96 100
61 126.3 0.0 271.0
98 59.3 0.0 130.2





#33

1,2-Dichloroethane-d4

Concen: 58.840 ug/l

RT: 5.940 min Scan# 796

Delta R.T. -0.000 min

Lab File: VX048735.D

Acq: 05 Dec 2025 14:54

Instrument :

MSVOA_X

ClientSampleId :

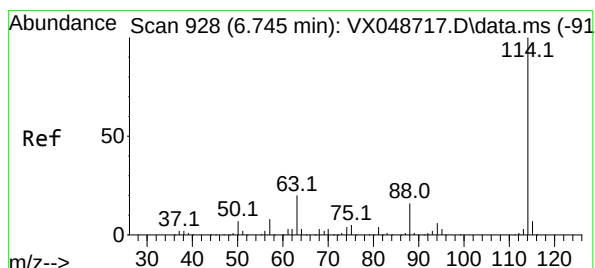
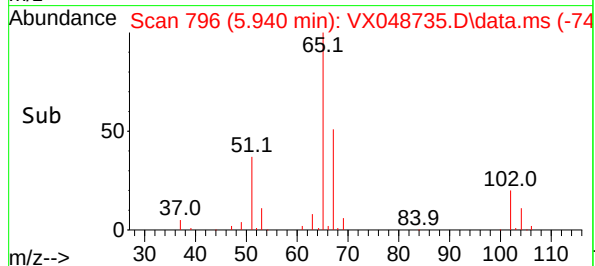
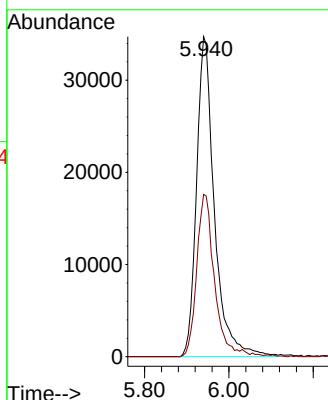
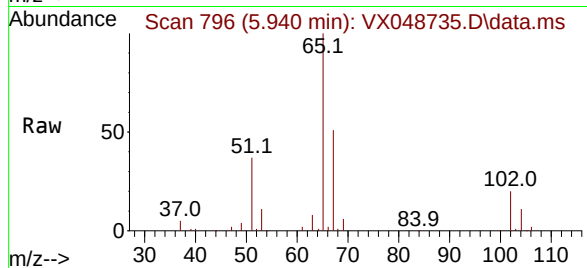
OW-05B-72-120325

Tgt Ion: 65 Resp: 102254

Ion Ratio Lower Upper

65 100

67 50.2 0.0 107.4



#34

1,4-Difluorobenzene

Concen: 50.000 ug/l

RT: 6.745 min Scan# 928

Delta R.T. -0.000 min

Lab File: VX048735.D

Acq: 05 Dec 2025 14:54

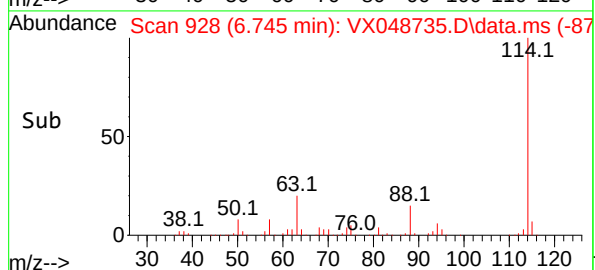
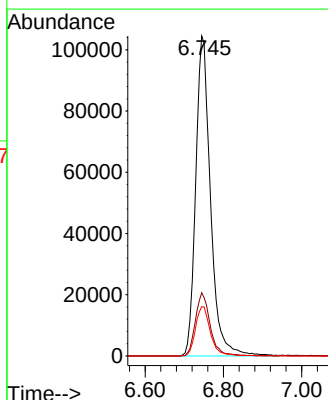
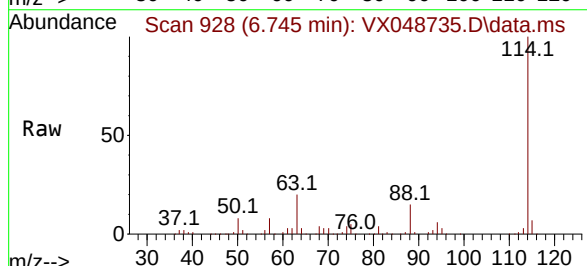
Tgt Ion: 114 Resp: 275147

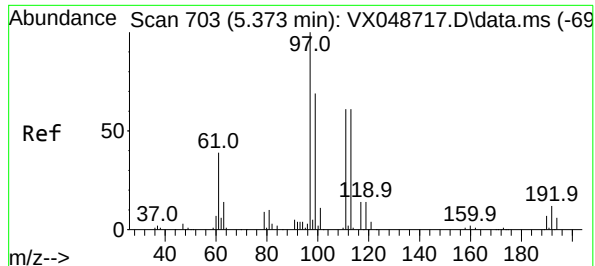
Ion Ratio Lower Upper

114 100

63 19.8 0.0 39.0

88 15.4 0.0 31.8





#35

Dibromofluoromethane

Concen: 39.688 ug/l

RT: 5.373 min Scan# 703

Delta R.T. -0.000 min

Lab File: VX048735.D

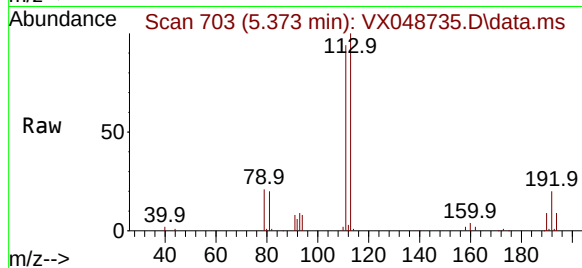
Acq: 05 Dec 2025 14:54

Instrument :

MSVOA_X

ClientSampleId :

OW-05B-72-120325



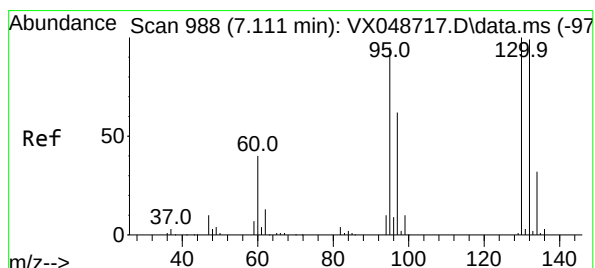
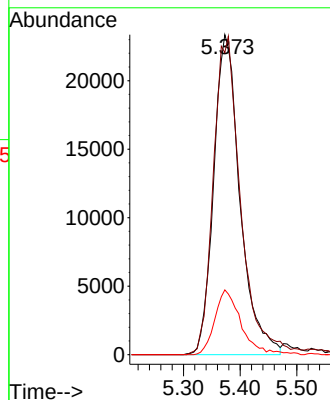
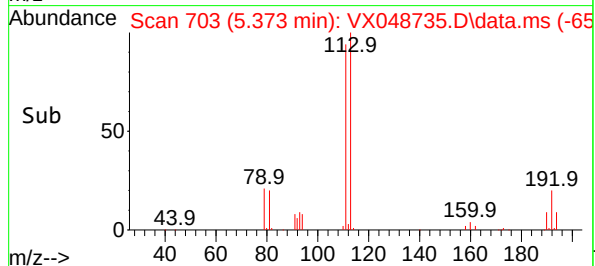
Tgt Ion:113 Resp: 75188

Ion Ratio Lower Upper

113 100

111 103.3 79.5 119.3

192 20.3 16.1 24.1



#44

Trichloroethene

Concen: 0.440 ug/l

RT: 7.110 min Scan# 988

Delta R.T. -0.000 min

Lab File: VX048735.D

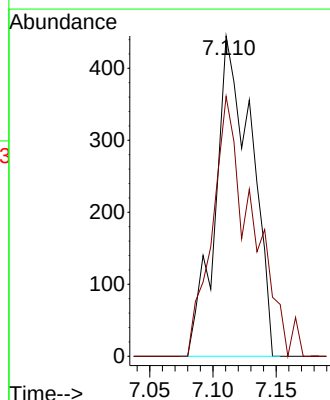
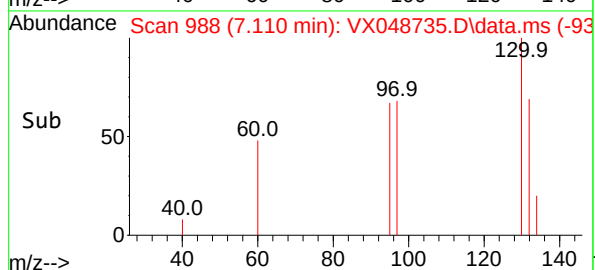
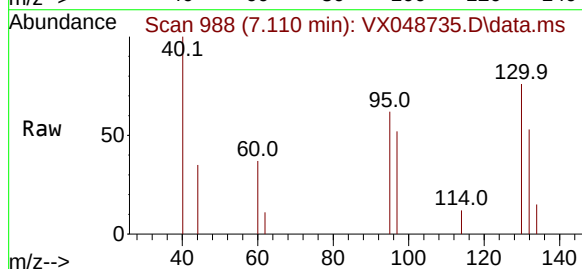
Acq: 05 Dec 2025 14:54

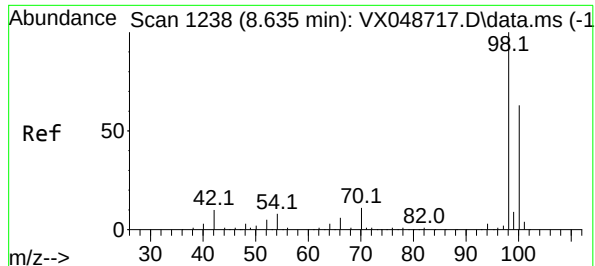
Tgt Ion:130 Resp: 883

Ion Ratio Lower Upper

130 100

95 81.1 0.0 189.6





#50

Toluene-d8

Concen: 38.182 ug/l

RT: 8.635 min Scan# 1

Delta R.T. -0.000 min

Lab File: VX048735.D

Acq: 05 Dec 2025 14:54

Instrument :

MSVOA_X

ClientSampleId :

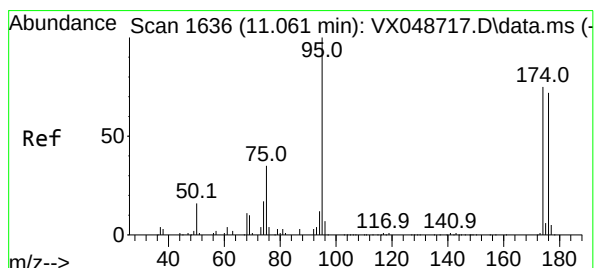
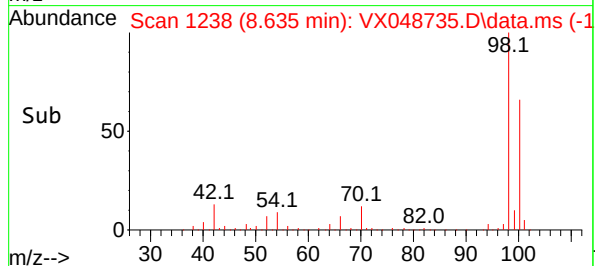
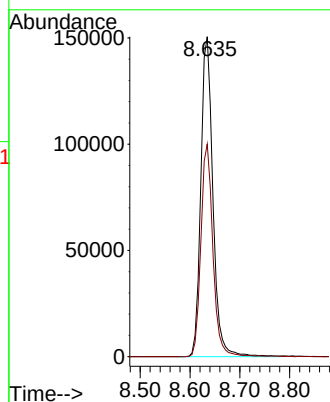
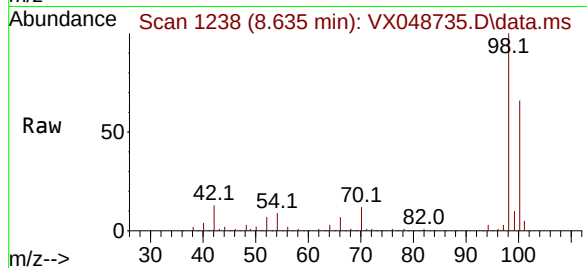
OW-05B-72-120325

Tgt Ion: 98 Resp: 253556

Ion Ratio Lower Upper

98 100

100 66.0 53.4 80.0



#62

4-Bromofluorobenzene

Concen: 45.048 ug/l

RT: 11.061 min Scan# 1636

Delta R.T. -0.000 min

Lab File: VX048735.D

Acq: 05 Dec 2025 14:54

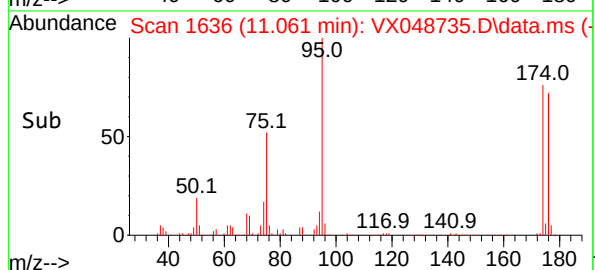
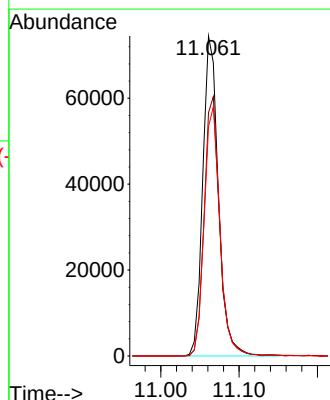
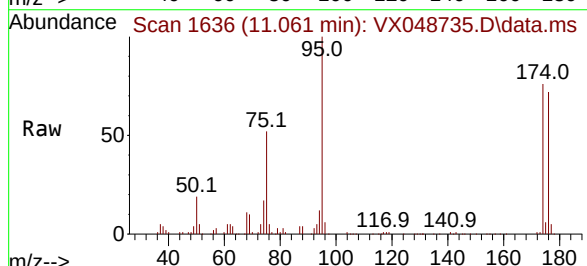
Tgt Ion: 95 Resp: 103159

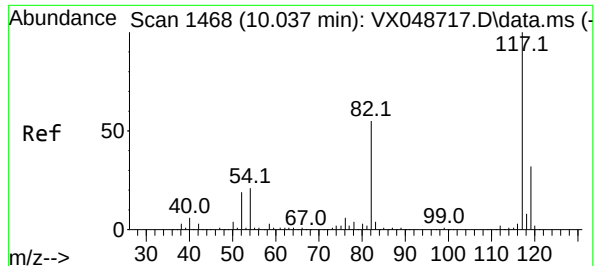
Ion Ratio Lower Upper

95 100

174 80.5 0.0 157.8

176 78.5 0.0 154.0





#63

Chlorobenzene-d5

Concen: 50.000 ug/l

RT: 10.037 min Scan# 1468

Delta R.T. -0.000 min

Lab File: VX048735.D

Acq: 05 Dec 2025 14:54

Instrument :

MSVOA_X

ClientSampleId :

OW-05B-72-120325

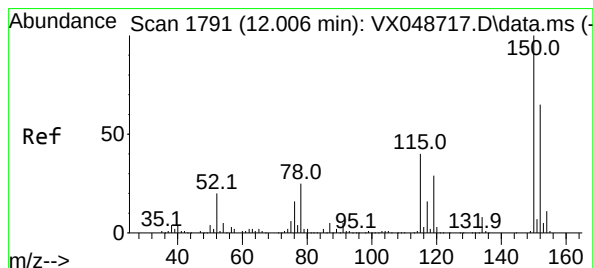
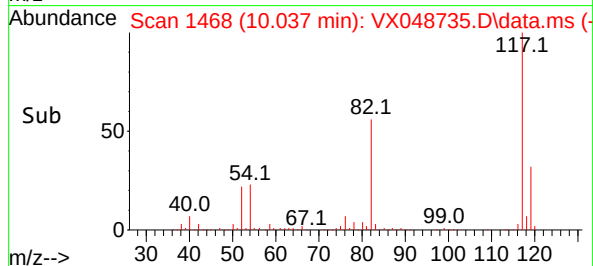
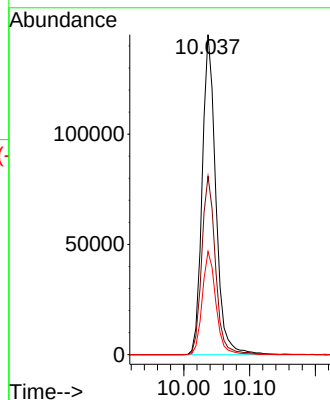
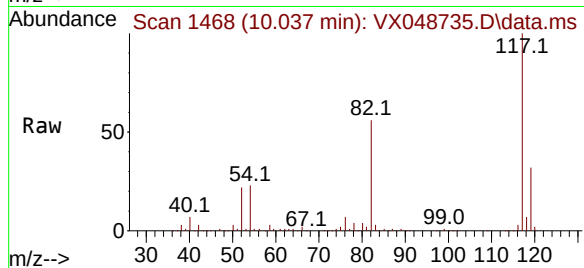
Tgt Ion:117 Resp: 208806

Ion Ratio Lower Upper

117 100

82 55.8 44.1 66.1

119 32.3 25.2 37.8



#72

1,4-Dichlorobenzene-d4

Concen: 50.000 ug/l

RT: 12.006 min Scan# 1791

Delta R.T. -0.000 min

Lab File: VX048735.D

Acq: 05 Dec 2025 14:54

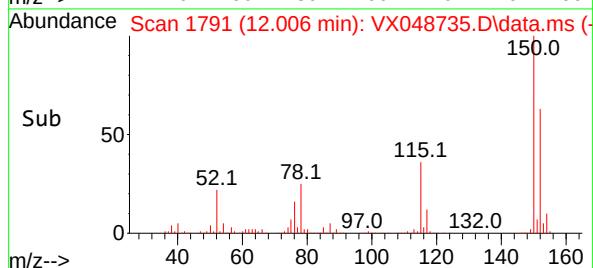
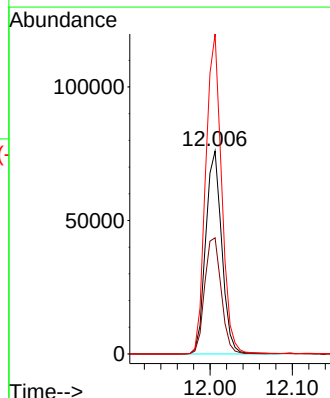
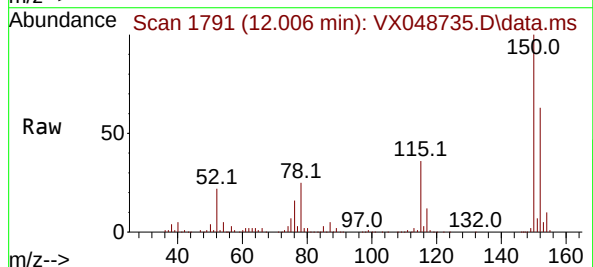
Tgt Ion:152 Resp: 103213

Ion Ratio Lower Upper

152 100

115 59.1 42.1 126.4

150 156.2 0.0 347.8





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

Report of Analysis

Client:	JACOBS Engineering Group, Inc.	Date Collected:	12/03/25
Project:	Former Schlumberger STC PTC Site D3868221	Date Received:	12/03/25
Client Sample ID:	OW-05B-72-120325RE	SDG No.:	Q3757
Lab Sample ID:	Q3757-09RE	Matrix:	Water
Analytical Method:	8260D	% Solid:	0
Sample Wt/Vol:	5 mL	Test:	VOCMS Group3
	Level: LOW		
	Final Vol: 5000 uL		

CAS Number	Parameter	Conc.	Qua.	DF	MDL	LOQ / CRQL	Units	Date Ana.	BatchID
TARGETS									
75-01-4	Vinyl Chloride	0.26	U	1	0.26	1.00	ug/L	12/08/25 12:53	VX120825
75-35-4	1,1-Dichloroethene	0.60	J	1	0.23	1.00	ug/L	12/08/25 12:53	VX120825
75-34-3	1,1-Dichloroethane	0.23	U	1	0.23	1.00	ug/L	12/08/25 12:53	VX120825
156-59-2	cis-1,2-Dichloroethene	1.50		1	0.19	1.00	ug/L	12/08/25 12:53	VX120825
71-55-6	1,1,1-Trichloroethane	0.20	U	1	0.20	1.00	ug/L	12/08/25 12:53	VX120825
71-43-2	Benzene	0.15	U	1	0.15	1.00	ug/L	12/08/25 12:53	VX120825
107-06-2	1,2-Dichloroethane	0.22	U	1	0.22	1.00	ug/L	12/08/25 12:53	VX120825
79-01-6	Trichloroethene	0.25	J	1	0.090	1.00	ug/L	12/08/25 12:53	VX120825
79-00-5	1,1,2-Trichloroethane	0.21	U	1	0.21	1.00	ug/L	12/08/25 12:53	VX120825
127-18-4	Tetrachloroethene	0.23	U	1	0.23	1.00	ug/L	12/08/25 12:53	VX120825
SURROGATES									
17060-07-0	1,2-Dichloroethane-d4	62.9			70 (74) - 130 (125)	126%	SPK: 50		
1868-53-7	Dibromofluoromethane	39.6			70 (75) - 130 (124)	79%	SPK: 50		
2037-26-5	Toluene-d8	47.0			70 (86) - 130 (113)	94%	SPK: 50		
460-00-4	4-Bromofluorobenzene	56.1			70 (77) - 130 (121)	112%	SPK: 50		
INTERNAL STANDARDS									
		Area Count							
363-72-4	Pentafluorobenzene	169000							
540-36-3	1,4-Difluorobenzene	410000							
3114-55-4	Chlorobenzene-d5	414000							
3855-82-1	1,4-Dichlorobenzene-d4	198000							

U = Not Detected

LOQ = Limit of Quantitation

MDL = Method Detection Limit

LOD = Limit of Detection

E = Value Exceeds Calibration Range

Q = indicates LCS control criteria did not meet requirements

M = MS/MSD acceptance criteria did not meet requirements

J = Estimated Value

B = Analyte Found in Associated Method Blank

N = Presumptive Evidence of a Compound

* = Values outside of QC limits

D = Dilution

() = Laboratory InHouse Limit

A = Aldol-Condensation Reaction Products

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120825\
 Data File : VX048748.D
 Acq On : 08 Dec 2025 12:53
 Operator : JC/MD
 Sample : Q3757-09RE
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 OW-05B-72-120325RE

Quant Time: Dec 09 04:09:02 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X120425W.M
 Quant Title : SW846 8260
 QLast Update : Fri Dec 05 03:27:34 2025
 Response via : Initial Calibration

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

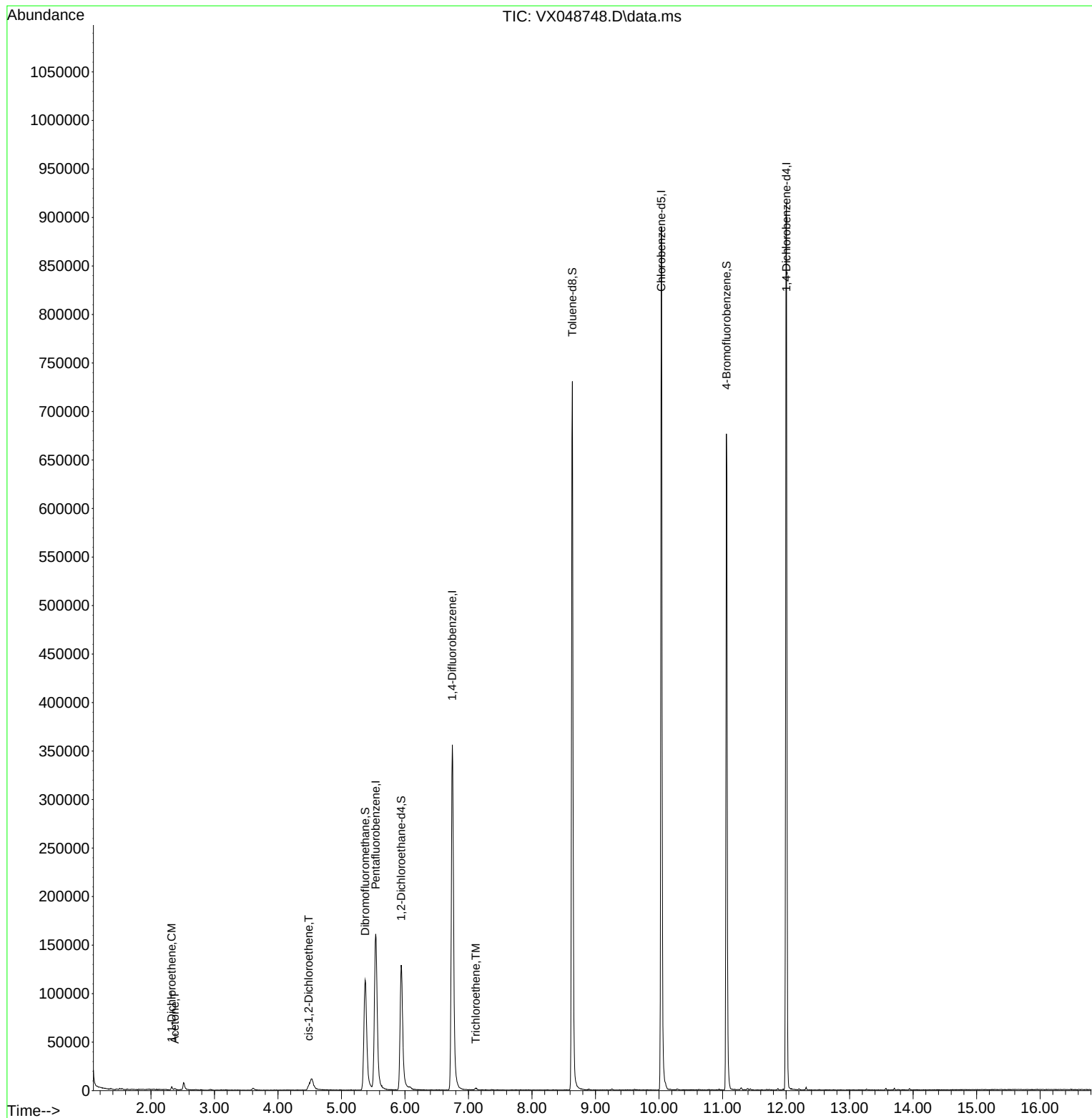
Internal Standards						
1) Pentafluorobenzene	5.537	168	169111	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.744	114	409831	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.036	117	413547	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.006	152	197728	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.940	65	142724	62.880	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	125.760%#
35) Dibromofluoromethane	5.373	113	111782	39.613	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	79.220%
50) Toluene-d8	8.634	98	465280	47.039	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	94.080%
62) 4-Bromofluorobenzene	11.061	95	191359	56.102	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	112.200%
Target Compounds						Qvalue
12) 1,1-Dichloroethene	2.324	96	1182	0.602	ug/l	86
16) Acetone	2.373	43	2059	2.370	ug/l	89
27) cis-1,2-Dichloroethene	4.483	96	3429	1.516	ug/l	60
44) Trichloroethene	7.116	130	749	0.251	ug/l	62

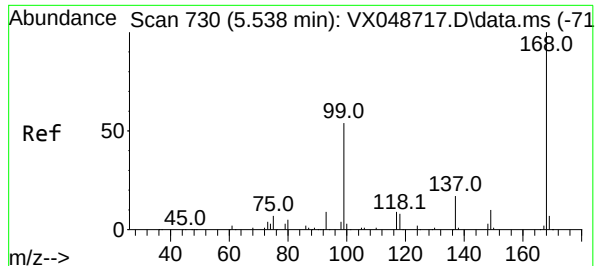
(#) = qualifier out of range (m) = manual integration (+) = signals summed

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120825\
Data File : VX048748.D
Acq On : 08 Dec 2025 12:53
Operator : JC/MD
Sample : Q3757-09RE
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
OW-05B-72-120325RE

Quant Time: Dec 09 04:09:02 2025
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X120425W.M
Quant Title : SW846 8260
QLast Update : Fri Dec 05 03:27:34 2025
Response via : Initial Calibration

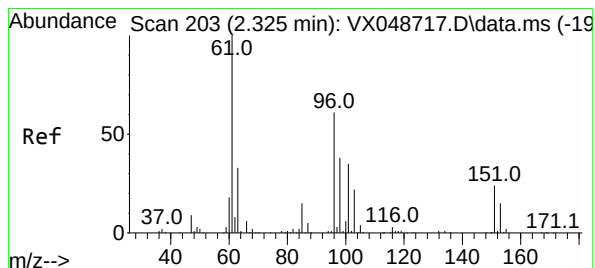
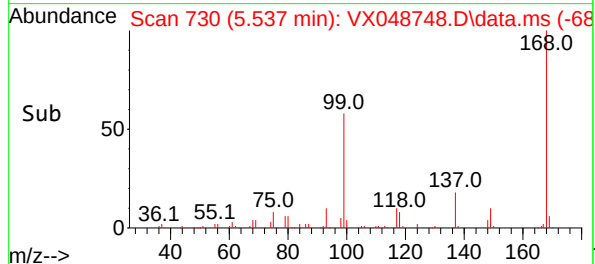
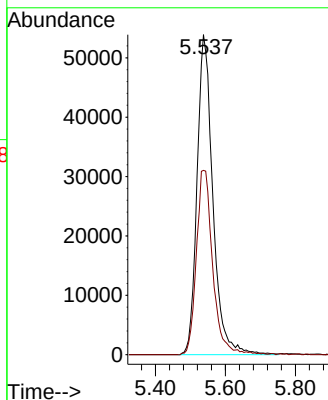
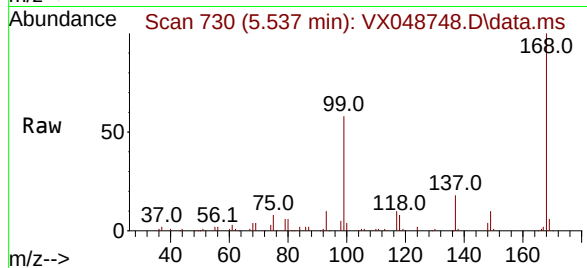




#1
Pentafluorobenzene
Concen: 50.000 ug/l
RT: 5.537 min Scan# 71
Delta R.T. -0.001 min
Lab File: VX048748.D
Acq: 08 Dec 2025 12:53

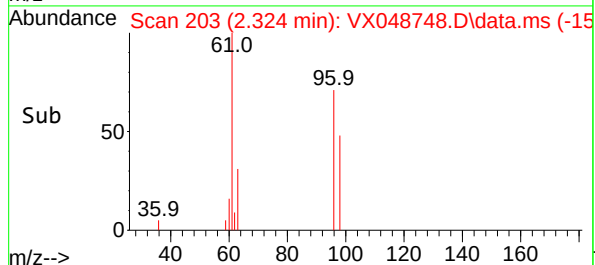
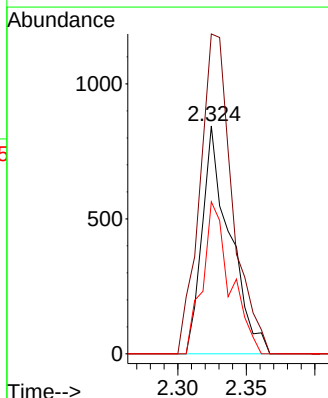
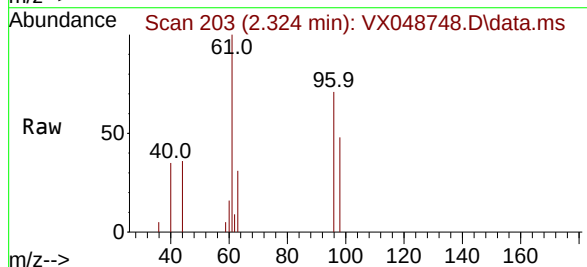
Instrument :
MSVOA_X
ClientSampleId :
OW-05B-72-120325RE

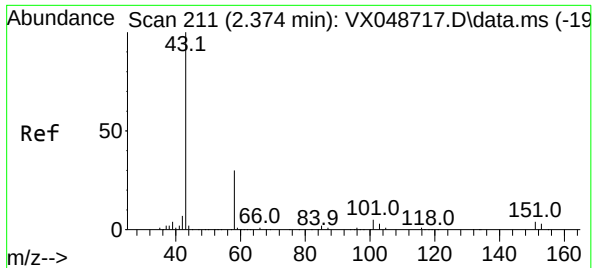
Tgt Ion:168 Resp: 169111
Ion Ratio Lower Upper
168 100
99 57.6 44.2 66.4



#12
1,1-Dichloroethene
Concen: 0.602 ug/l
RT: 2.324 min Scan# 203
Delta R.T. -0.000 min
Lab File: VX048748.D
Acq: 08 Dec 2025 12:53

Tgt Ion: 96 Resp: 1182
Ion Ratio Lower Upper
96 100
61 140.4 130.7 196.1
98 66.7 49.8 74.8

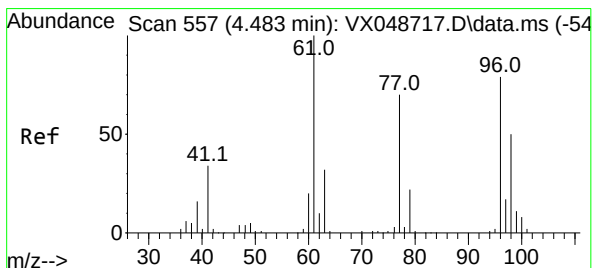
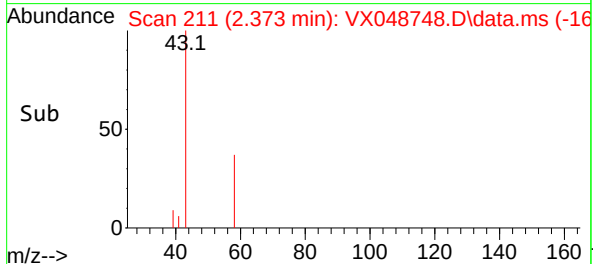
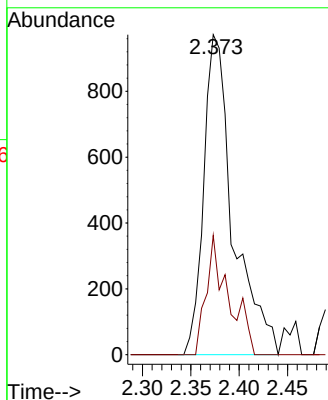
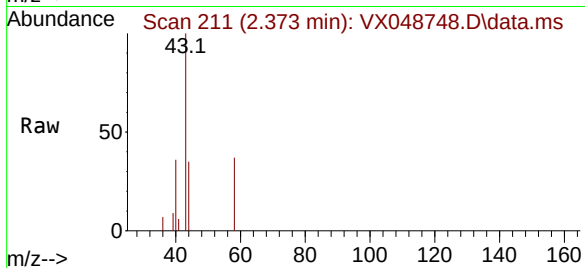




#16
Acetone
Concen: 2.370 ug/l
RT: 2.373 min Scan# 211
Delta R.T. -0.000 min
Lab File: VX048748.D
Acq: 08 Dec 2025 12:53

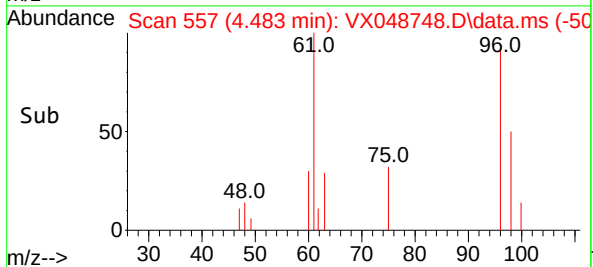
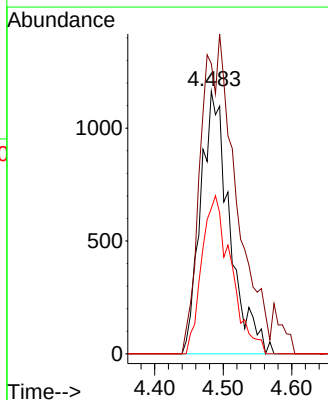
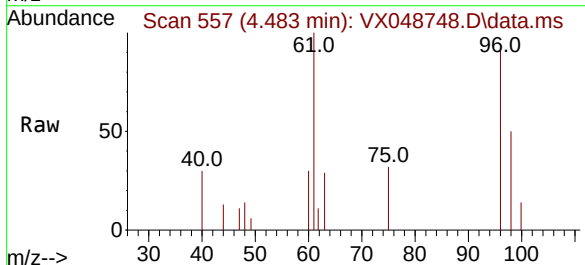
Instrument :
MSVOA_X
ClientSampleId :
OW-05B-72-120325RE

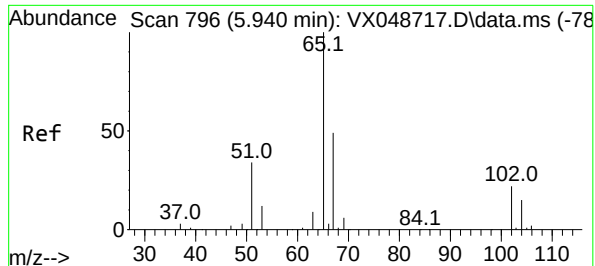
Tgt Ion: 43 Resp: 2059
Ion Ratio Lower Upper
43 100
58 37.2 25.0 37.4



#27
cis-1,2-Dichloroethene
Concen: 1.516 ug/l
RT: 4.483 min Scan# 557
Delta R.T. -0.000 min
Lab File: VX048748.D
Acq: 08 Dec 2025 12:53

Tgt Ion: 96 Resp: 3429
Ion Ratio Lower Upper
96 100
61 67.6 0.0 271.0
98 61.1 0.0 130.2





#33

1,2-Dichloroethane-d4

Concen: 62.880 ug/l

RT: 5.940 min Scan# 796

Delta R.T. -0.000 min

Lab File: VX048748.D

Acq: 08 Dec 2025 12:53

Instrument :

MSVOA_X

ClientSampleId :

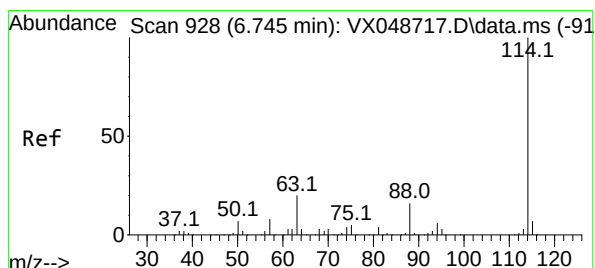
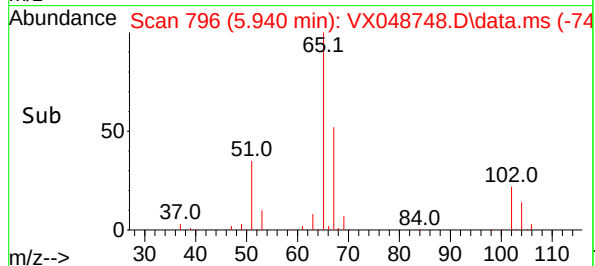
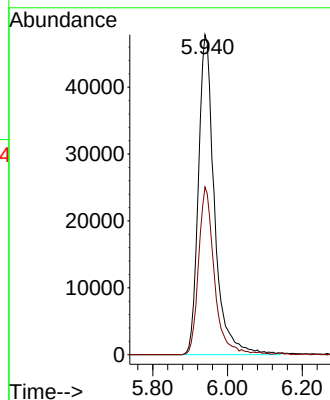
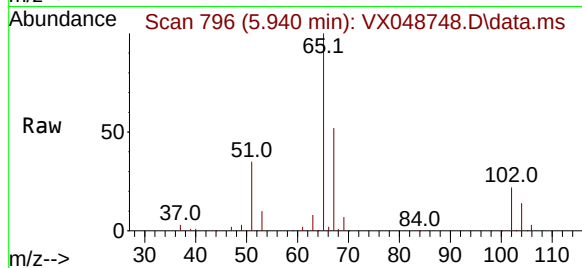
OW-05B-72-120325RE

Tgt Ion: 65 Resp: 142724

Ion Ratio Lower Upper

65 100

67 51.7 0.0 107.4



#34

1,4-Difluorobenzene

Concen: 50.000 ug/l

RT: 6.744 min Scan# 928

Delta R.T. -0.000 min

Lab File: VX048748.D

Acq: 08 Dec 2025 12:53

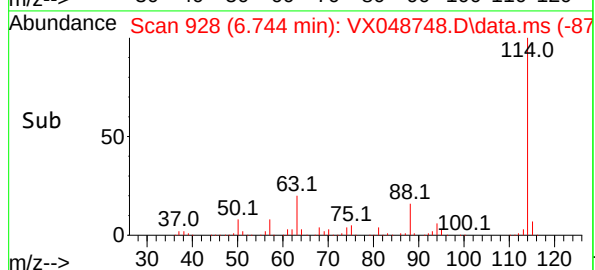
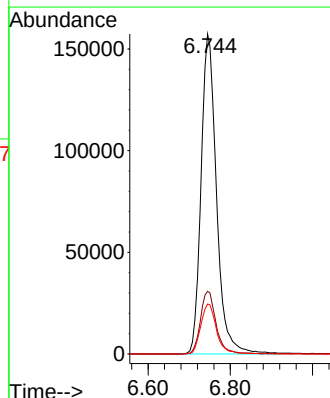
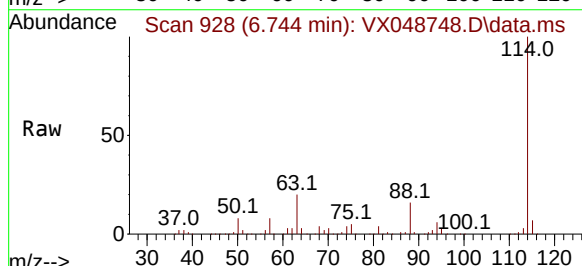
Tgt Ion: 114 Resp: 409831

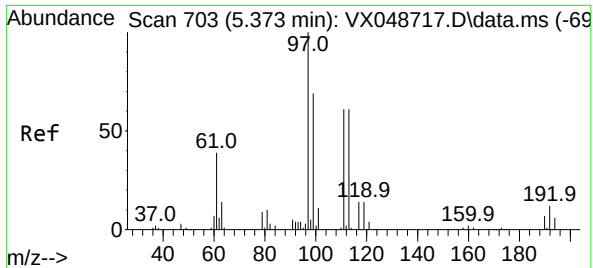
Ion Ratio Lower Upper

114 100

63 19.5 0.0 39.0

88 15.6 0.0 31.8





#35

Dibromofluoromethane

Concen: 39.613 ug/l

RT: 5.373 min Scan# 703

Delta R.T. -0.000 min

Lab File: VX048748.D

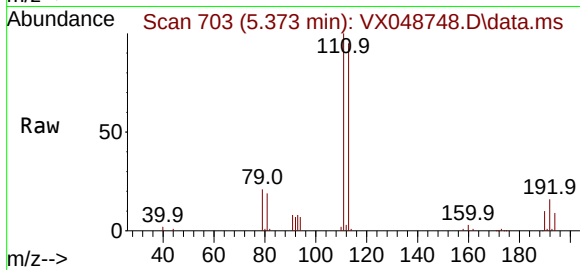
Acq: 08 Dec 2025 12:53

Instrument :

MSVOA_X

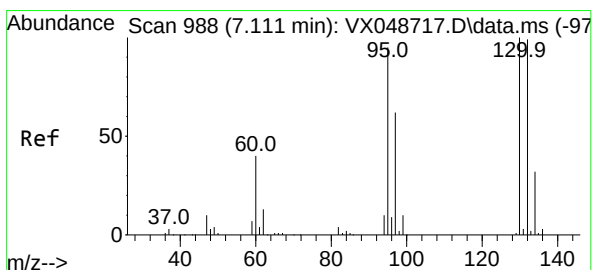
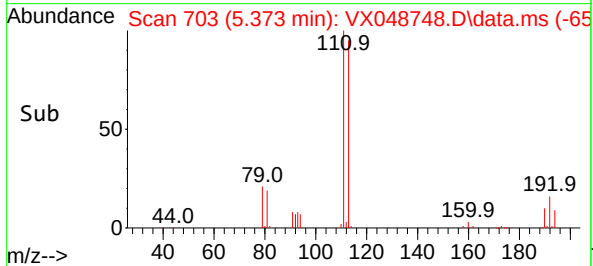
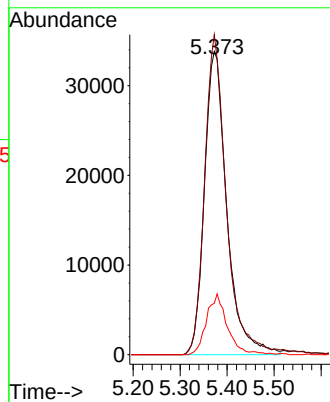
ClientSampleId :

OW-05B-72-120325RE



Tgt Ion:113 Resp: 111782

Ion	Ratio	Lower	Upper
113	100		
111	103.7	79.5	119.3
192	18.7	16.1	24.1



#44

Trichloroethene

Concen: 0.251 ug/l

RT: 7.116 min Scan# 989

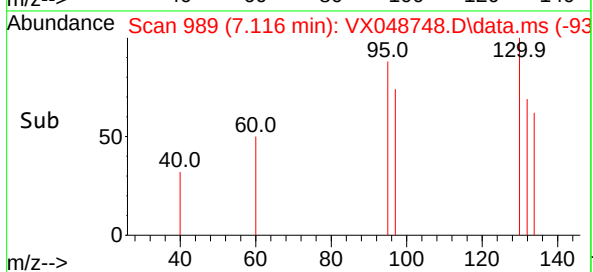
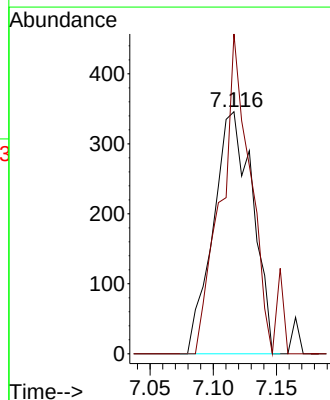
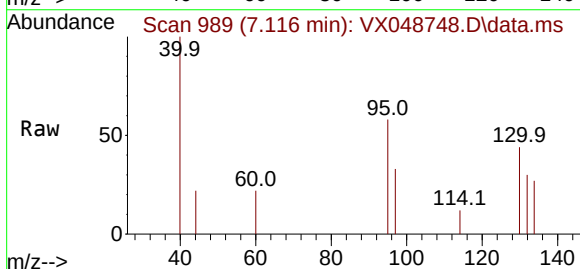
Delta R.T. 0.006 min

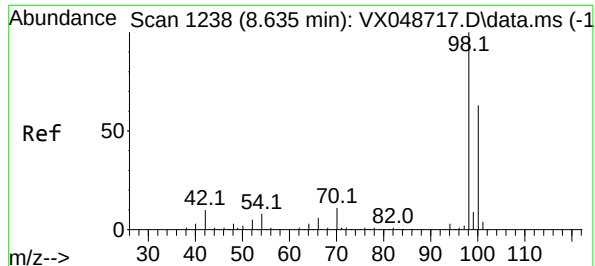
Lab File: VX048748.D

Acq: 08 Dec 2025 12:53

Tgt Ion:130 Resp: 749

Ion	Ratio	Lower	Upper
130	100		
95	132.1	0.0	189.6





#50

Toluene-d8

Concen: 47.039 ug/l

RT: 8.634 min Scan# 1

Delta R.T. -0.000 min

Lab File: VX048748.D

Acq: 08 Dec 2025 12:53

Instrument :

MSVOA_X

ClientSampleId :

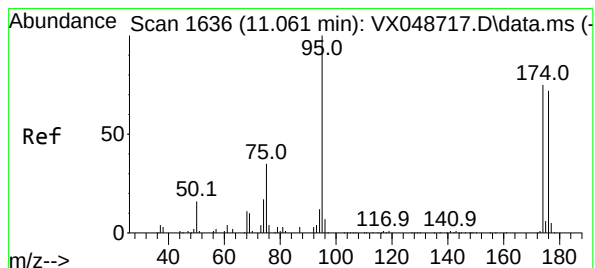
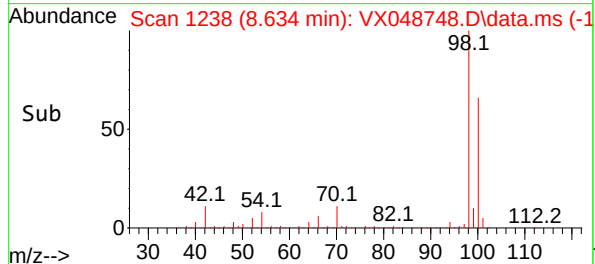
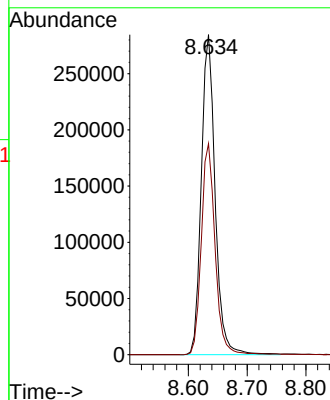
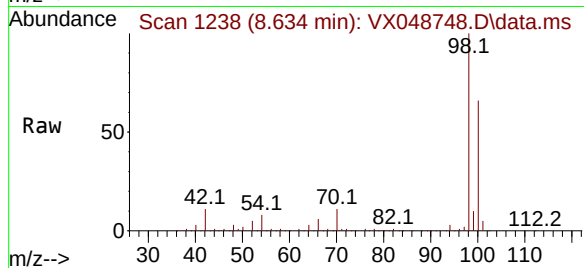
OW-05B-72-120325RE

Tgt Ion: 98 Resp: 465280

Ion Ratio Lower Upper

98 100

100 66.2 53.4 80.0



#62

4-Bromofluorobenzene

Concen: 56.102 ug/l

RT: 11.061 min Scan# 1636

Delta R.T. -0.000 min

Lab File: VX048748.D

Acq: 08 Dec 2025 12:53

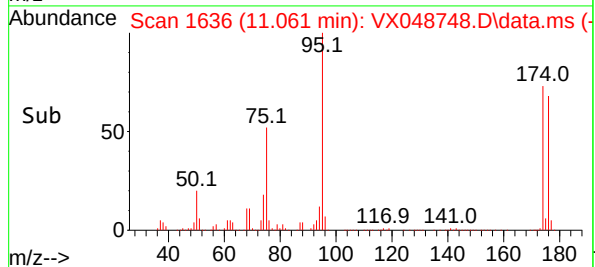
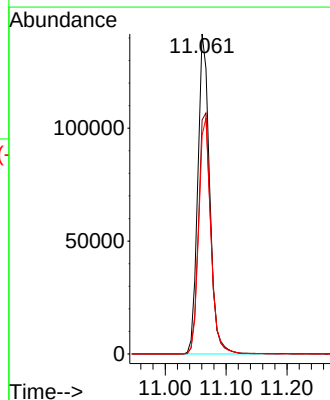
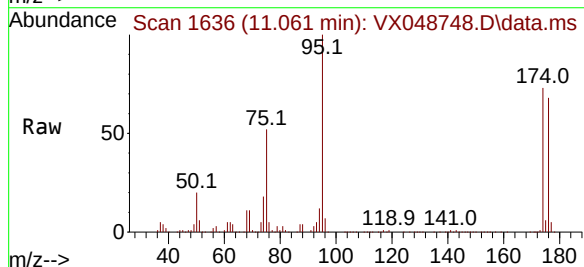
Tgt Ion: 95 Resp: 191359

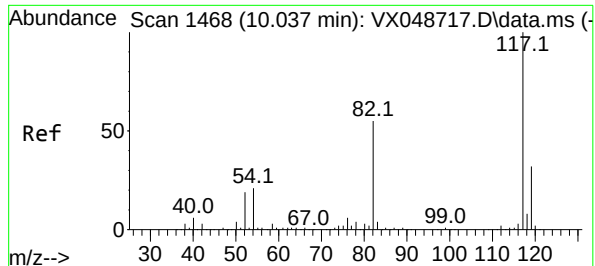
Ion Ratio Lower Upper

95 100

174 77.9 0.0 157.8

176 75.0 0.0 154.0





#63

Chlorobenzene-d5

Concen: 50.000 ug/l

RT: 10.036 min Scan# 1468

Delta R.T. -0.000 min

Lab File: VX048748.D

Acq: 08 Dec 2025 12:53

Instrument :

MSVOA_X

ClientSampleId :

OW-05B-72-120325RE

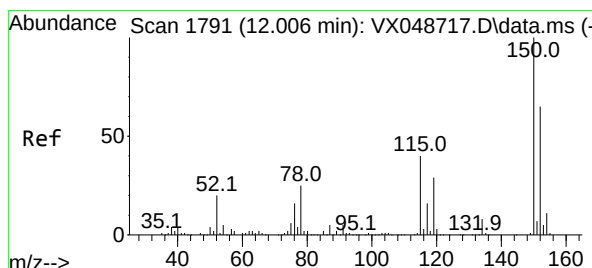
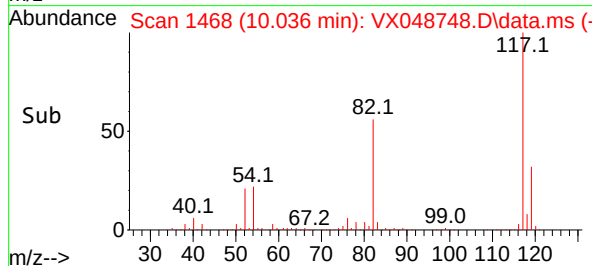
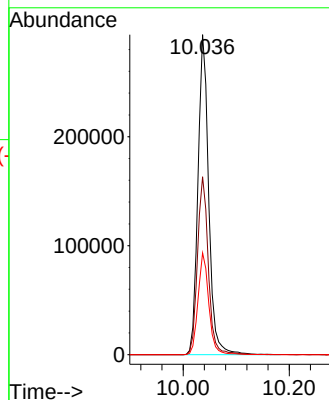
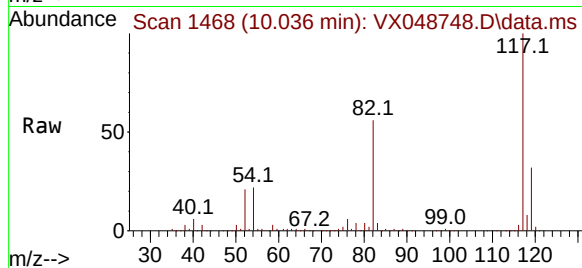
Tgt Ion:117 Resp: 413547

Ion Ratio Lower Upper

117 100

82 55.6 44.1 66.1

119 31.8 25.2 37.8



#72

1,4-Dichlorobenzene-d4

Concen: 50.000 ug/l

RT: 12.006 min Scan# 1791

Delta R.T. -0.000 min

Lab File: VX048748.D

Acq: 08 Dec 2025 12:53

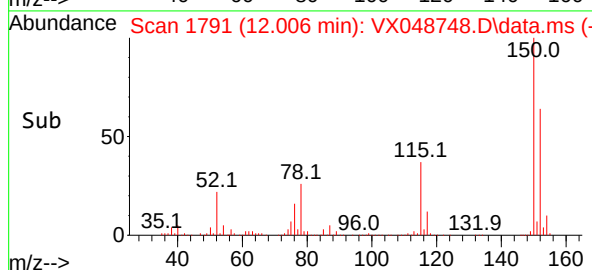
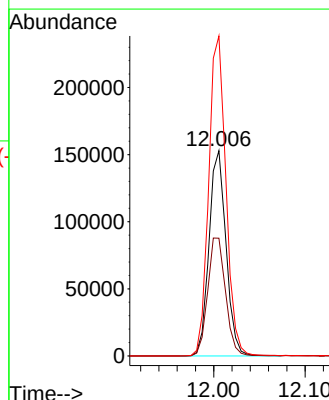
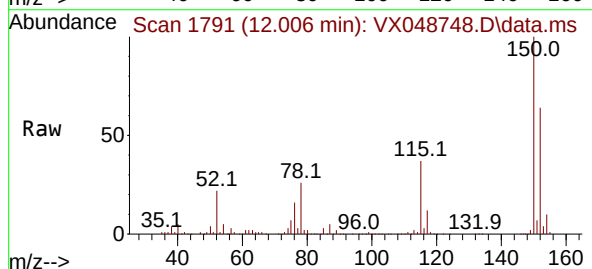
Tgt Ion:152 Resp: 197728

Ion Ratio Lower Upper

152 100

115 60.2 42.1 126.4

150 157.5 0.0 347.8





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

Report of Analysis

Client: JACOBS Engineering Group, Inc.
Project: Former Schlumberger STC PTC Site D3868221
Client Sample ID: EB01-120325
Lab Sample ID: Q3757-11
Analytical Method: 8260D
Sample Wt/Vol: 5 mL

Level: LOW
Final Vol: 5000 uL

Date Collected: 12/03/25
Date Received: 12/03/25
SDG No.: Q3757
Matrix: Water
% Solid: 0
Test: VOCMS Group3

CAS Number	Parameter	Conc.	Qua.	DF	MDL	LOQ / CRQL	Units	Date Ana.	BatchID
TARGETS									
75-01-4	Vinyl Chloride	0.26	U	1	0.26	1.00	ug/L	12/04/25 16:48	VN120425
75-35-4	1,1-Dichloroethene	0.23	U	1	0.23	1.00	ug/L	12/04/25 16:48	VN120425
75-34-3	1,1-Dichloroethane	0.23	U	1	0.23	1.00	ug/L	12/04/25 16:48	VN120425
156-59-2	cis-1,2-Dichloroethene	0.19	U	1	0.19	1.00	ug/L	12/04/25 16:48	VN120425
71-55-6	1,1,1-Trichloroethane	0.20	U	1	0.20	1.00	ug/L	12/04/25 16:48	VN120425
71-43-2	Benzene	0.15	U	1	0.15	1.00	ug/L	12/04/25 16:48	VN120425
107-06-2	1,2-Dichloroethane	0.22	U	1	0.22	1.00	ug/L	12/04/25 16:48	VN120425
79-01-6	Trichloroethene	0.090	U	1	0.090	1.00	ug/L	12/04/25 16:48	VN120425
79-00-5	1,1,2-Trichloroethane	0.21	U	1	0.21	1.00	ug/L	12/04/25 16:48	VN120425
127-18-4	Tetrachloroethene	0.23	U	1	0.23	1.00	ug/L	12/04/25 16:48	VN120425
SURROGATES									
17060-07-0	1,2-Dichloroethane-d4	53.6			70 (74) - 130 (125)	107%	SPK: 50		
1868-53-7	Dibromofluoromethane	46.9			70 (75) - 130 (124)	94%	SPK: 50		
2037-26-5	Toluene-d8	44.3			70 (86) - 130 (113)	89%	SPK: 50		
460-00-4	4-Bromofluorobenzene	49.5			70 (77) - 130 (121)	99%	SPK: 50		
INTERNAL STANDARDS									
		Area Count							
363-72-4	Pentafluorobenzene	202000							
540-36-3	1,4-Difluorobenzene	452000							
3114-55-4	Chlorobenzene-d5	440000							
3855-82-1	1,4-Dichlorobenzene-d4	211000							

U = Not Detected

LOQ = Limit of Quantitation

MDL = Method Detection Limit

LOD = Limit of Detection

E = Value Exceeds Calibration Range

Q = indicates LCS control criteria did not meet requirements

M = MS/MSD acceptance criteria did not meet requirements

J = Estimated Value

B = Analyte Found in Associated Method Blank

N = Presumptive Evidence of a Compound

* = Values outside of QC limits

D = Dilution

() = Laboratory InHouse Limit

A = Aldol-Condensation Reaction Products

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN120425\
 Data File : VN088450.D
 Acq On : 04 Dec 2025 16:48
 Operator : JC\MD
 Sample : Q3757-11
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 EB01-120325

Manual Integrations
 APPROVED

Reviewed By :John Carlone 12/05/2025
 Supervised By :Semsettin Yesilyurt 12/05/2025

Quant Time: Dec 05 00:18:07 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N112425W.M
 Quant Title : SW846 8260
 QLast Update : Tue Nov 25 01:34:52 2025
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.206	168	201753	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.082	114	451562	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.841	117	439722	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.770	152	210586	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.559	65	203838	53.633	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery = 107.260%			
35) Dibromofluoromethane	8.147	113	146660	46.867	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery = 93.740%			
50) Toluene-d8	10.541	98	516378	44.349	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery = 88.700%			
62) 4-Bromofluorobenzene	12.829	95	197862	49.526	ug/l	0.00
Spiked Amount 50.000	Range 77 - 121		Recovery = 99.060%			
Target Compounds					Qvalue	
16) Acetone	4.412	43	17456m	12.700	ug/l	

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

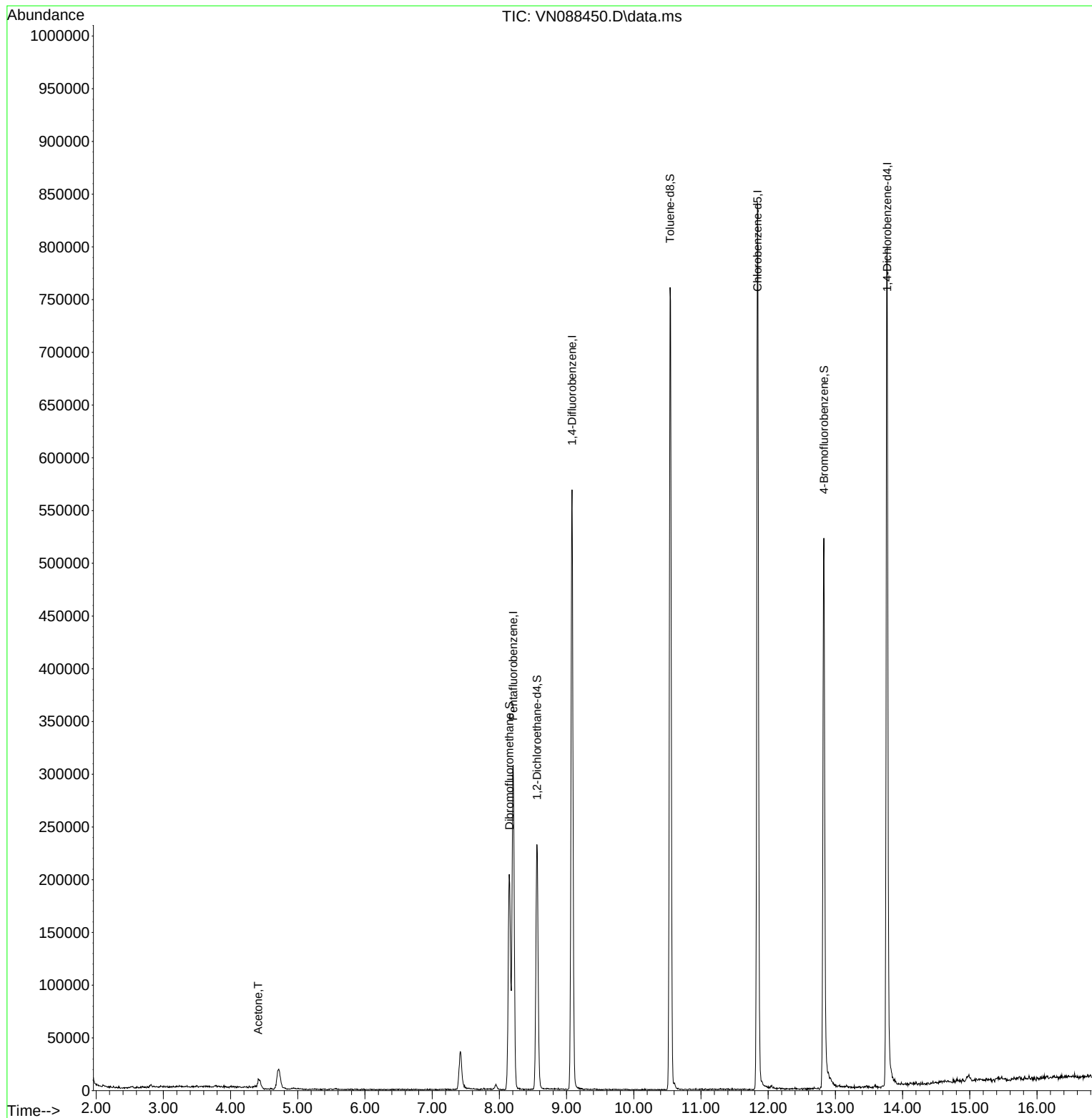
Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN120425\
Data File : VN088450.D
Acq On : 04 Dec 2025 16:48
Operator : JC\MD
Sample : Q3757-11
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 16 Sample Multiplier: 1

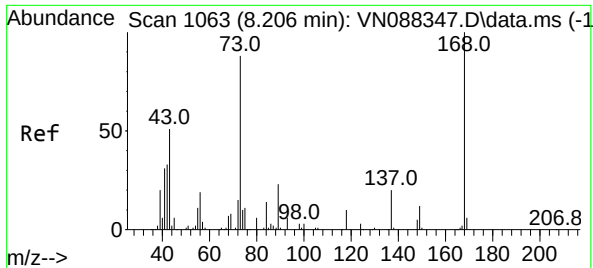
Instrument :
MSVOA_N
ClientSampleId :
EB01-120325

Manual Integrations
APPROVED

Reviewed By :John Carlone 12/05/2025
Supervised By :Semsettin Yesilyurt 12/05/2025

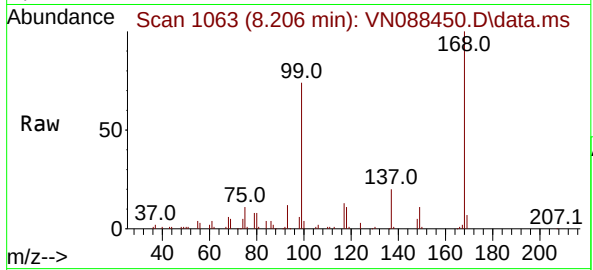
Quant Time: Dec 05 00:18:07 2025
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N112425W.M
Quant Title : SW846 8260
QLast Update : Tue Nov 25 01:34:52 2025
Response via : Initial Calibration





#1
Pentafluorobenzene
Concen: 50.000 ug/l
RT: 8.206 min Scan# 1063
Delta R.T. 0.000 min
Lab File: VN088450.D
Acq: 04 Dec 2025 16:48

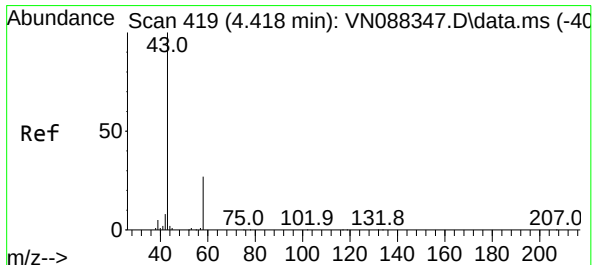
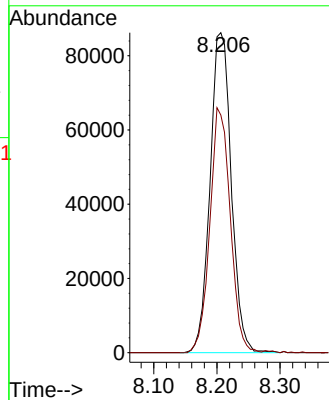
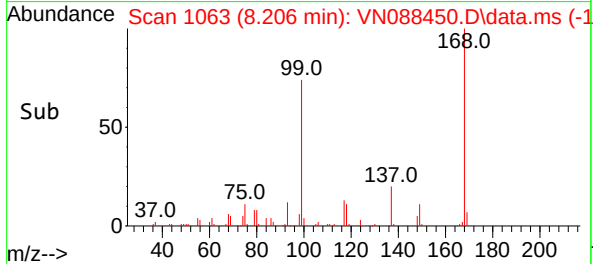
Instrument :
MSVOA_N
ClientSampleId :
EB01-120325



Tgt Ion:168 Resp: 20175
Ion Ratio Lower Upper
168 100
99 74.3 57.1 85.7

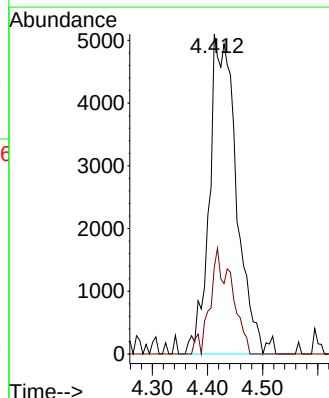
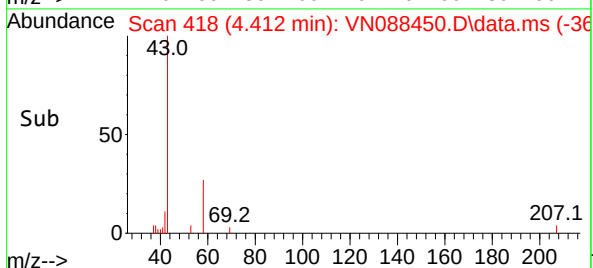
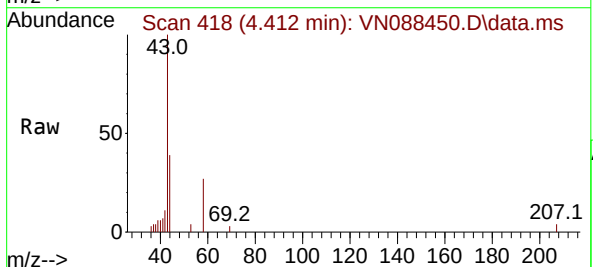
Manual Integrations
APPROVED

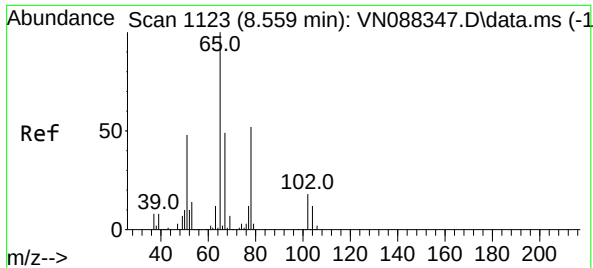
Reviewed By :John Carlone 12/05/2025
Supervised By :Semsettin Yesilyurt 12/05/2025



#16
Acetone
Concen: 12.700 ug/l m
RT: 4.412 min Scan# 418
Delta R.T. -0.006 min
Lab File: VN088450.D
Acq: 04 Dec 2025 16:48

Tgt Ion: 43 Resp: 17456
Ion Ratio Lower Upper
43 100
58 27.1 21.7 32.5





#33

1,2-Dichloroethane-d4

Concen: 53.633 ug/l

RT: 8.559 min Scan# 1123

Delta R.T. 0.000 min

Lab File: VN088450.D

Acq: 04 Dec 2025 16:48

Instrument :

MSVOA_N

ClientSampleId :

EB01-120325

Tgt Ion: 65 Resp: 203838

Ion Ratio Lower Upper

65 100

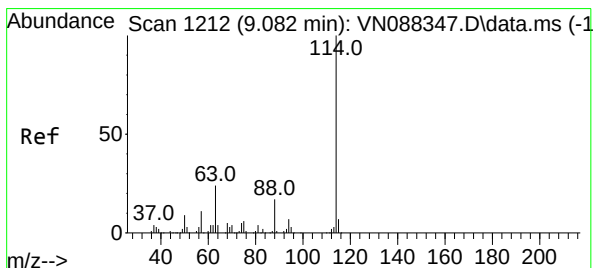
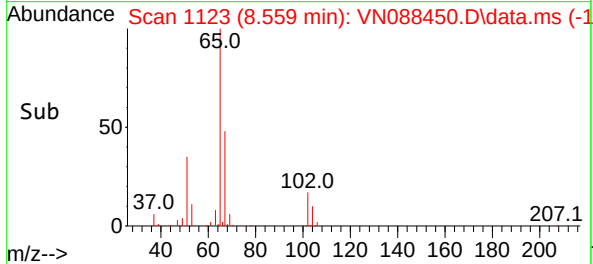
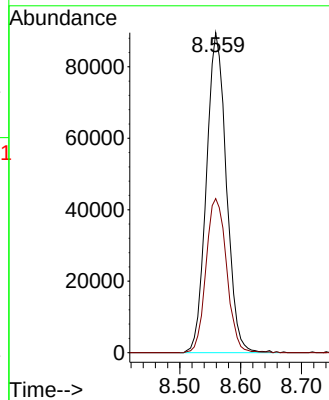
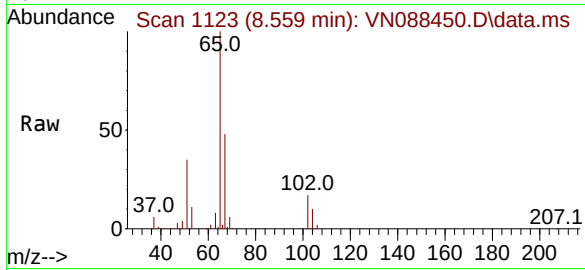
67 49.7 0.0 100.8

Manual Integrations

APPROVED

Reviewed By :John Carlone 12/05/2025

Supervised By :Semsettin Yesilyurt 12/05/2025



#34

1,4-Difluorobenzene

Concen: 50.000 ug/l

RT: 9.082 min Scan# 1212

Delta R.T. 0.000 min

Lab File: VN088450.D

Acq: 04 Dec 2025 16:48

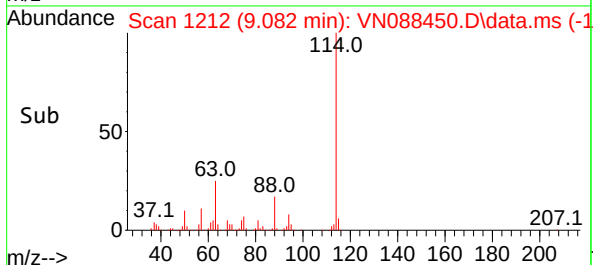
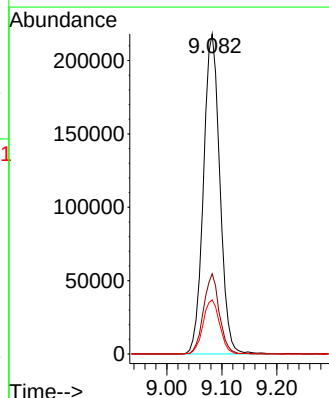
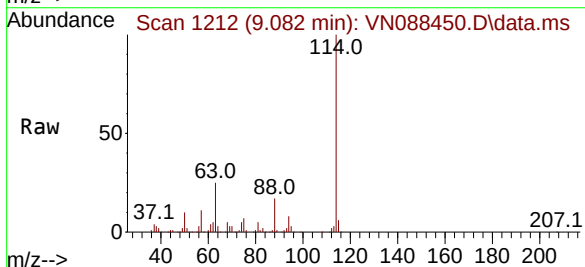
Tgt Ion:114 Resp: 451562

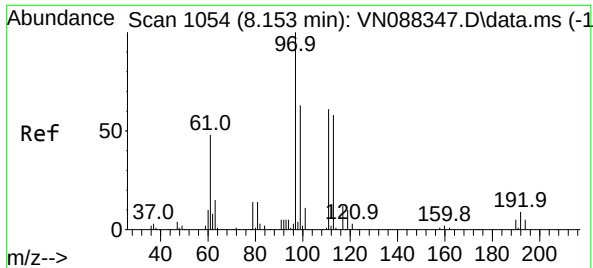
Ion Ratio Lower Upper

114 100

63 25.0 0.0 49.0

88 16.9 0.0 34.0





#35

Dibromofluoromethane

Concen: 46.867 ug/l

RT: 8.147 min Scan# 1053

Delta R.T. -0.006 min

Lab File: VN088450.D

Acq: 04 Dec 2025 16:48

Instrument :

MSVOA_N

ClientSampleId :

EB01-120325

Tgt Ion:113 Resp: 146660

Ion Ratio Lower Upper

113 100

111 101.9 82.5 123.7

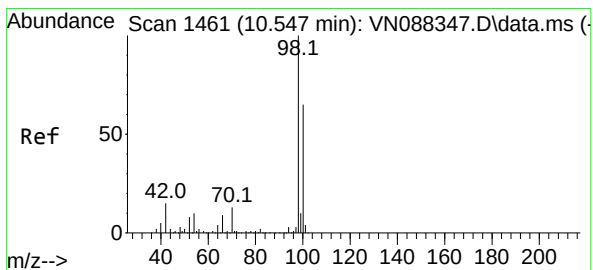
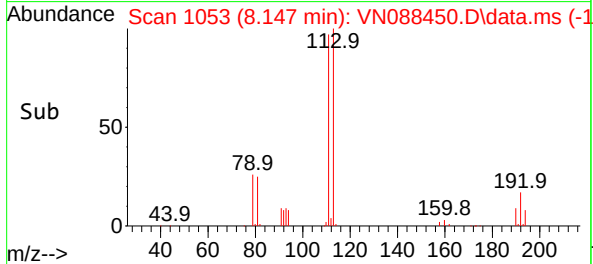
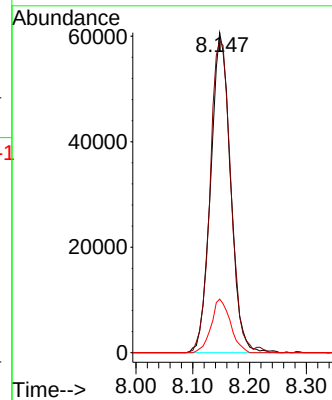
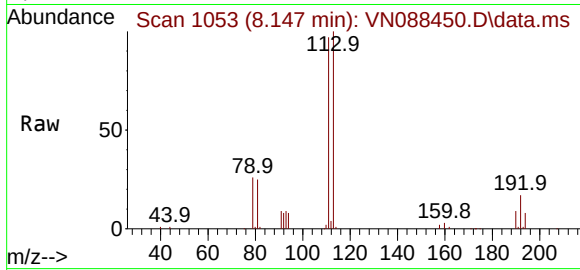
192 16.2 12.8 19.2

Manual Integrations

APPROVED

Reviewed By :John Carlone 12/05/2025

Supervised By :Semsettin Yesilyurt 12/05/2025



#50

Toluene-d8

Concen: 44.349 ug/l

RT: 10.541 min Scan# 1460

Delta R.T. -0.006 min

Lab File: VN088450.D

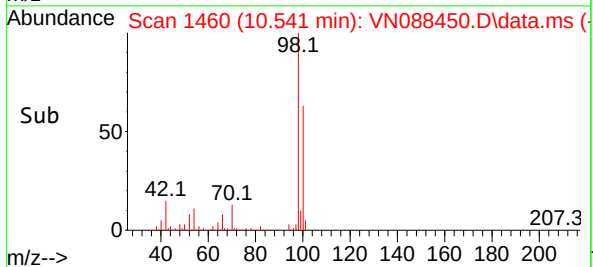
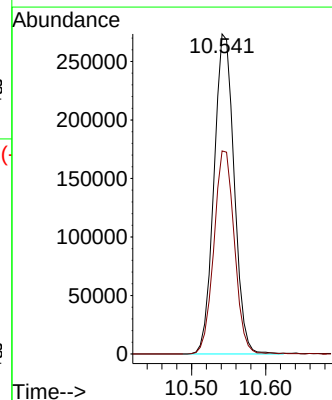
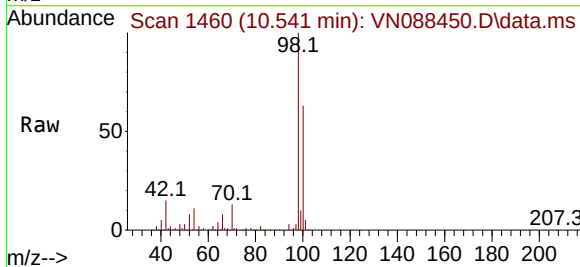
Acq: 04 Dec 2025 16:48

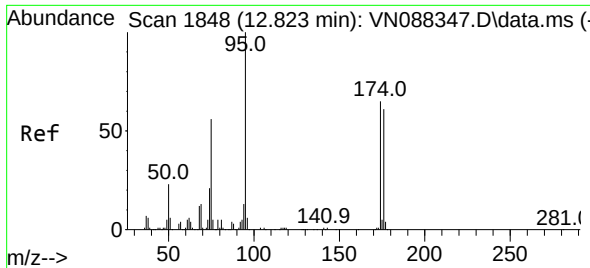
Tgt Ion: 98 Resp: 516378

Ion Ratio Lower Upper

98 100

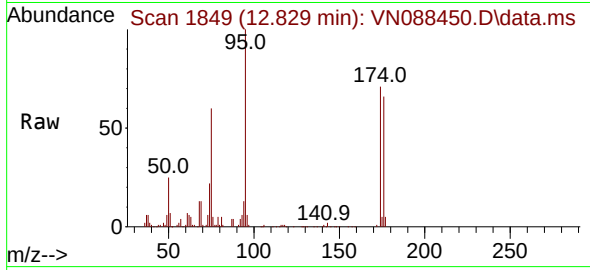
100 64.5 52.2 78.2





#62
4-Bromofluorobenzene
Concen: 49.526 ug/l
RT: 12.829 min Scan# 1848
Delta R.T. 0.006 min
Lab File: VN088450.D
Acq: 04 Dec 2025 16:48

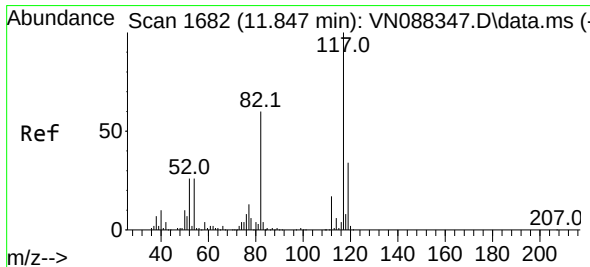
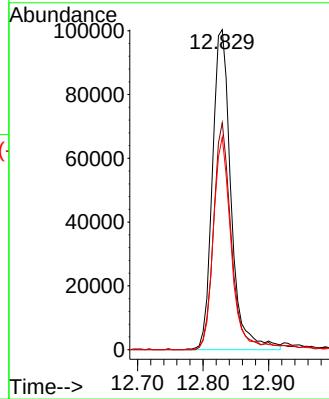
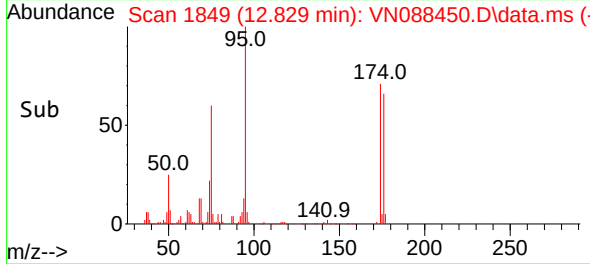
Instrument :
MSVOA_N
ClientSampleId :
EB01-120325



Tgt Ion: 95 Resp: 197862
Ion Ratio Lower Upper
95 100
174 66.5 0.0 139.0
176 64.3 0.0 132.4

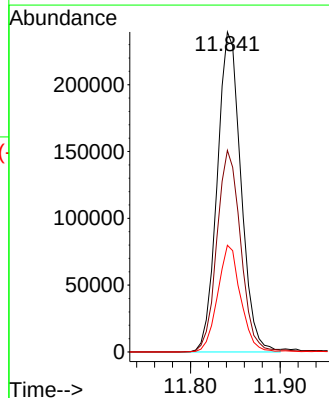
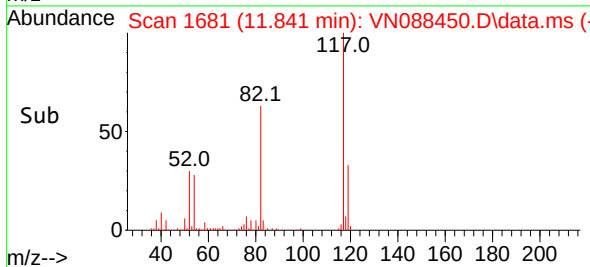
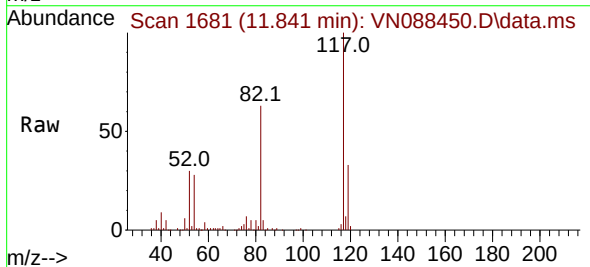
Manual Integrations
APPROVED

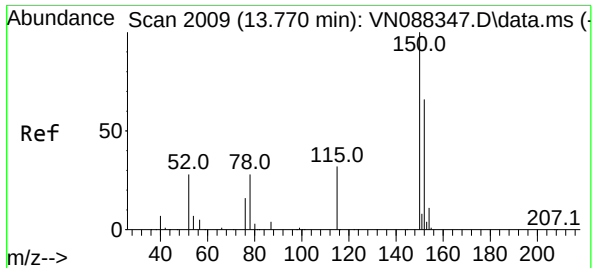
Reviewed By :John Carlone 12/05/2025
Supervised By :Semsettin Yesilyurt 12/05/2025



#63
Chlorobenzene-d5
Concen: 50.000 ug/l
RT: 11.841 min Scan# 1681
Delta R.T. -0.006 min
Lab File: VN088450.D
Acq: 04 Dec 2025 16:48

Tgt Ion:117 Resp: 439722
Ion Ratio Lower Upper
117 100
82 63.0 47.8 71.8
119 33.4 26.9 40.3





#72

1,4-Dichlorobenzene-d4

Concen: 50.000 ug/l

RT: 13.770 min Scan# 20

Delta R.T. 0.000 min

Lab File: VN088450.D

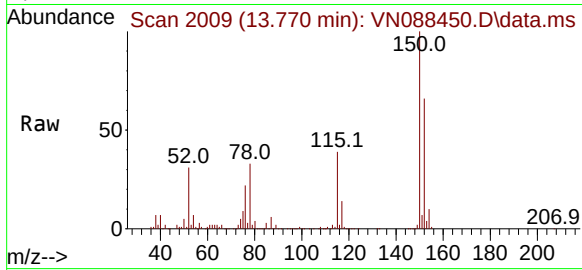
Acq: 04 Dec 2025 16:48

Instrument :

MSVOA_N

ClientSampleId :

EB01-120325



Tgt Ion:152 Resp: 210580

Ion Ratio Lower Upper

152 100

115 66.3 33.0 98.9

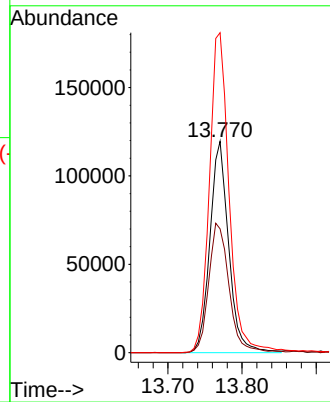
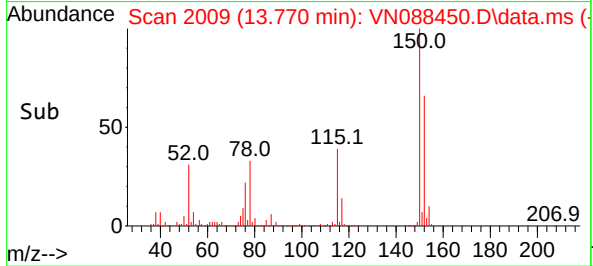
150 160.2 0.0 349.0

Manual Integrations

APPROVED

Reviewed By :John Carlone 12/05/2025

Supervised By :Semsettin Yesilyurt 12/05/2025





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

Report of Analysis

Client: JACOBS Engineering Group, Inc.
Project: Former Schlumberger STC PTC Site D3868221
Client Sample ID: TB01-120325
Lab Sample ID: Q3757-12
Analytical Method: 8260D
Sample Wt/Vol: 5 mL

Level: LOW
Final Vol: 5000 uL

Date Collected: 12/03/25
Date Received: 12/03/25
SDG No.: Q3757
Matrix: Water
% Solid: 0
Test: VOCMS Group3

CAS Number	Parameter	Conc.	Qua.	DF	MDL	LOQ / CRQL	Units	Date Ana.	BatchID
TARGETS									
75-01-4	Vinyl Chloride	0.26	U	1	0.26	1.00	ug/L	12/04/25 16:28	VN120425
75-35-4	1,1-Dichloroethene	0.23	U	1	0.23	1.00	ug/L	12/04/25 16:28	VN120425
75-34-3	1,1-Dichloroethane	0.23	U	1	0.23	1.00	ug/L	12/04/25 16:28	VN120425
156-59-2	cis-1,2-Dichloroethene	0.19	U	1	0.19	1.00	ug/L	12/04/25 16:28	VN120425
71-55-6	1,1,1-Trichloroethane	0.20	U	1	0.20	1.00	ug/L	12/04/25 16:28	VN120425
71-43-2	Benzene	0.15	U	1	0.15	1.00	ug/L	12/04/25 16:28	VN120425
107-06-2	1,2-Dichloroethane	0.22	U	1	0.22	1.00	ug/L	12/04/25 16:28	VN120425
79-01-6	Trichloroethene	0.090	U	1	0.090	1.00	ug/L	12/04/25 16:28	VN120425
79-00-5	1,1,2-Trichloroethane	0.21	U	1	0.21	1.00	ug/L	12/04/25 16:28	VN120425
127-18-4	Tetrachloroethene	0.23	U	1	0.23	1.00	ug/L	12/04/25 16:28	VN120425
SURROGATES									
17060-07-0	1,2-Dichloroethane-d4	56.4			70 (74) - 130 (125)	113%	SPK: 50		
1868-53-7	Dibromofluoromethane	47.1			70 (75) - 130 (124)	94%	SPK: 50		
2037-26-5	Toluene-d8	45.0			70 (86) - 130 (113)	90%	SPK: 50		
460-00-4	4-Bromofluorobenzene	52.4			70 (77) - 130 (121)	105%	SPK: 50		
INTERNAL STANDARDS									
		Area Count							
363-72-4	Pentafluorobenzene	198000							
540-36-3	1,4-Difluorobenzene	454000							
3114-55-4	Chlorobenzene-d5	441000							
3855-82-1	1,4-Dichlorobenzene-d4	211000							

U = Not Detected

LOQ = Limit of Quantitation

MDL = Method Detection Limit

LOD = Limit of Detection

E = Value Exceeds Calibration Range

Q = indicates LCS control criteria did not meet requirements

M = MS/MSD acceptance criteria did not meet requirements

J = Estimated Value

B = Analyte Found in Associated Method Blank

N = Presumptive Evidence of a Compound

* = Values outside of QC limits

D = Dilution

() = Laboratory InHouse Limit

A = Aldol-Condensation Reaction Products

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN120425\
Data File : VN088449.D
Acq On : 04 Dec 2025 16:28
Operator : JC\MD
Sample : Q3757-12
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
TB01-120325

Quant Time: Dec 05 00:17:51 2025
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N112425W.M
Quant Title : SW846 8260
QLast Update : Tue Nov 25 01:34:52 2025
Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.206	168	198228	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.083	114	454339	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.841	117	441078	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.770	152	211349	50.000	ug/l	0.00

System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.559	65	210629	56.405	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	112.820%
35) Dibromofluoromethane	8.147	113	148440	47.146	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	94.300%
50) Toluene-d8	10.541	98	527723	45.047	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	90.100%
62) 4-Bromofluorobenzene	12.829	95	210610	52.395	ug/l	0.00
Spiked Amount	50.000	Range	77 - 121	Recovery	=	104.800%

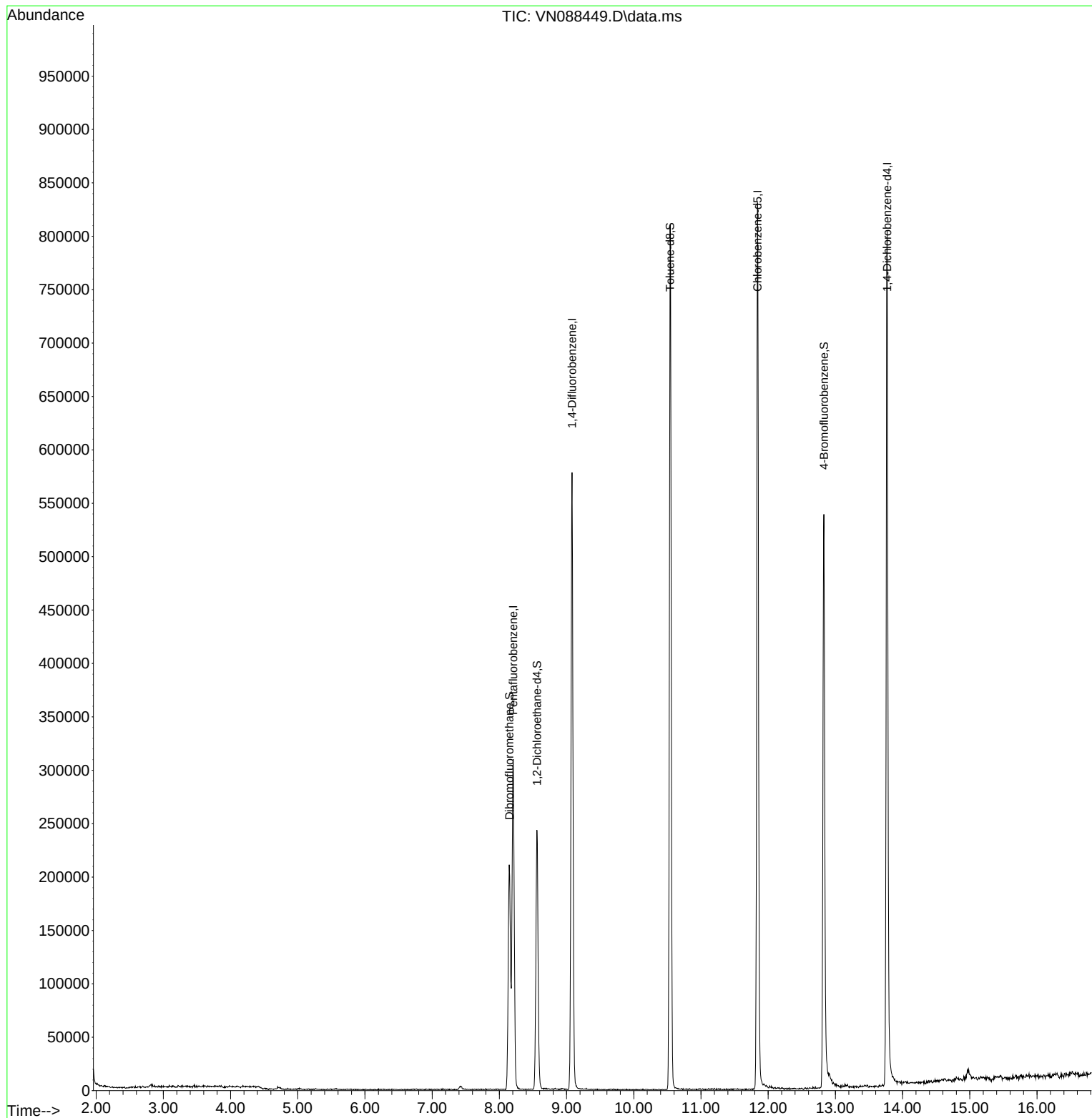
Target Compounds	Qvalue

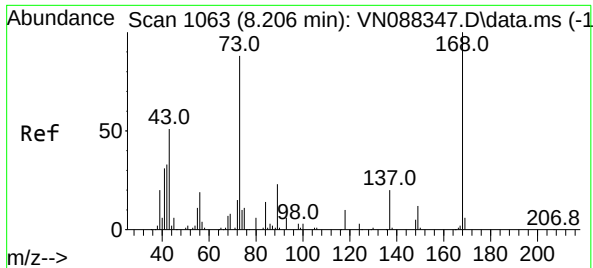
(#) = qualifier out of range (m) = manual integration (+) = signals summed

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN120425\
Data File : VN088449.D
Acq On : 04 Dec 2025 16:28
Operator : JC\MD
Sample : Q3757-12
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
TB01-120325

Quant Time: Dec 05 00:17:51 2025
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N112425W.M
Quant Title : SW846 8260
QLast Update : Tue Nov 25 01:34:52 2025
Response via : Initial Calibration

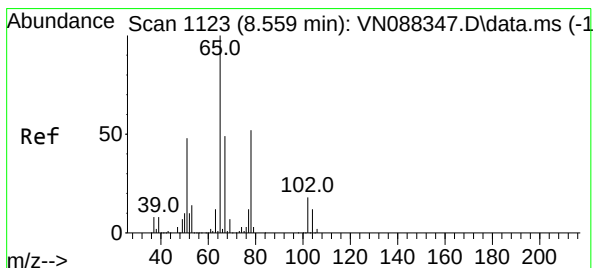
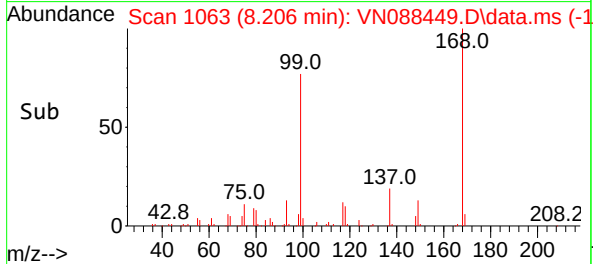
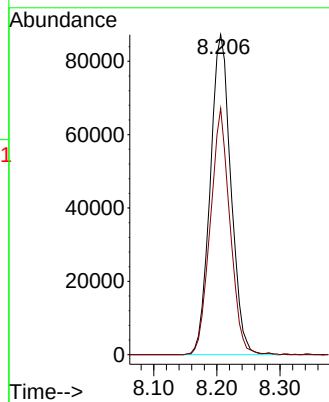
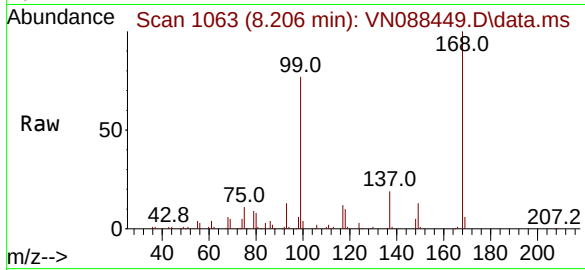




#1
Pentafluorobenzene
Concen: 50.000 ug/l
RT: 8.206 min Scan# 1063
Delta R.T. 0.000 min
Lab File: VN088449.D
Acq: 04 Dec 2025 16:28

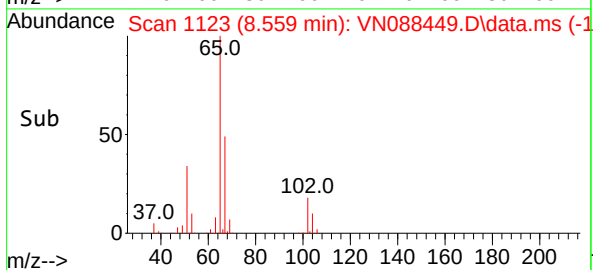
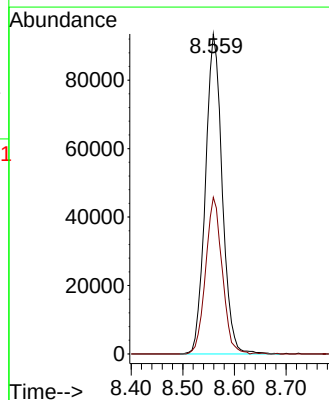
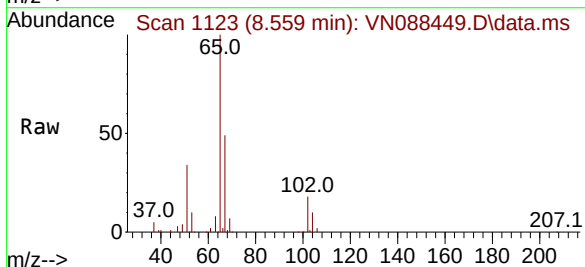
Instrument :
MSVOA_N
ClientSampleId :
TB01-120325

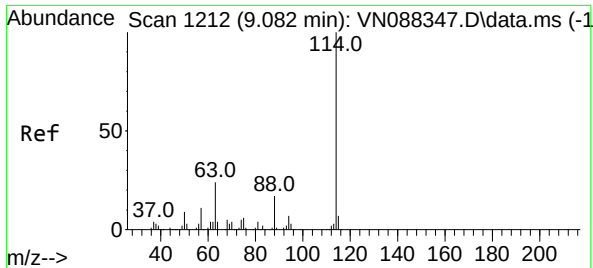
Tgt Ion:168 Resp: 198228
Ion Ratio Lower Upper
168 100
99 77.0 57.1 85.7



#33
1,2-Dichloroethane-d4
Concen: 56.405 ug/l
RT: 8.559 min Scan# 1123
Delta R.T. 0.000 min
Lab File: VN088449.D
Acq: 04 Dec 2025 16:28

Tgt Ion: 65 Resp: 210629
Ion Ratio Lower Upper
65 100
67 48.4 0.0 100.8





#34

1,4-Difluorobenzene

Concen: 50.000 ug/l

RT: 9.083 min Scan# 11

Delta R.T. 0.000 min

Lab File: VN088449.D

Acq: 04 Dec 2025 16:28

Instrument :

MSVOA_N

ClientSampleId :

TB01-120325

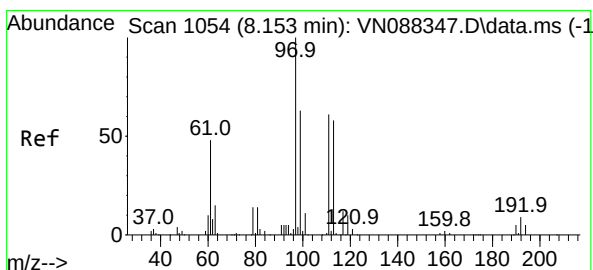
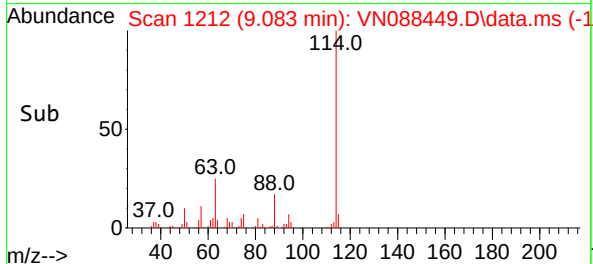
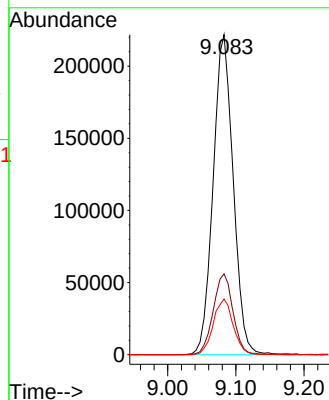
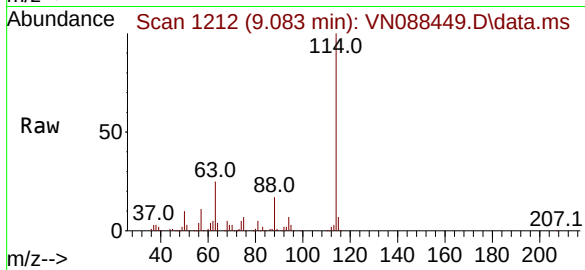
Tgt Ion:114 Resp: 454339

Ion Ratio Lower Upper

114 100

63 25.2 0.0 49.0

88 17.4 0.0 34.0



#35

Dibromofluoromethane

Concen: 47.146 ug/l

RT: 8.147 min Scan# 1053

Delta R.T. -0.006 min

Lab File: VN088449.D

Acq: 04 Dec 2025 16:28

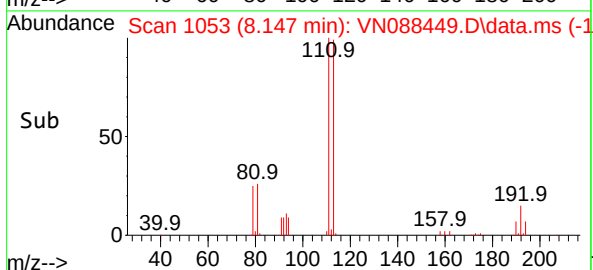
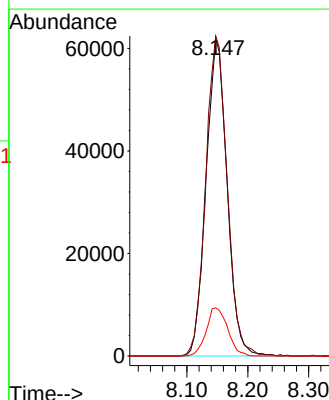
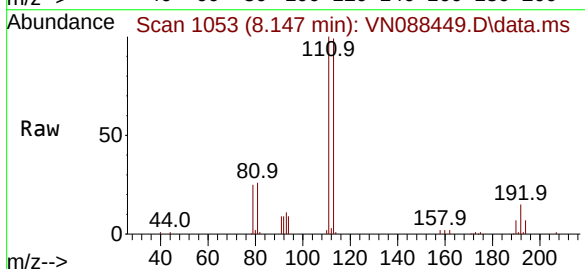
Tgt Ion:113 Resp: 148440

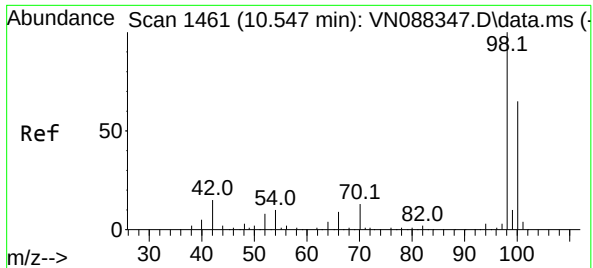
Ion Ratio Lower Upper

113 100

111 102.4 82.5 123.7

192 15.4 12.8 19.2

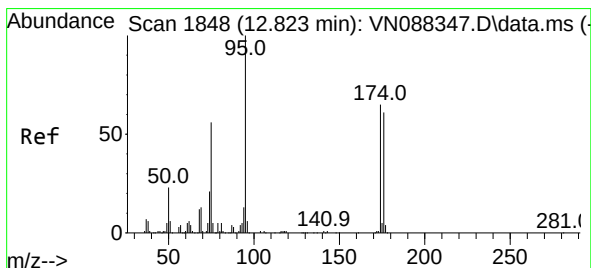
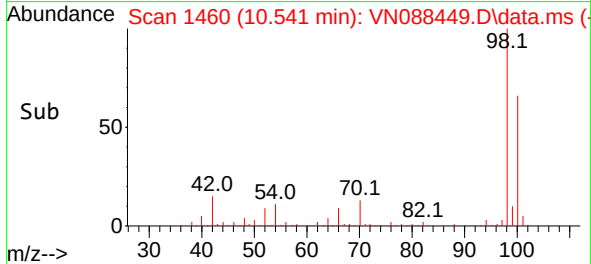
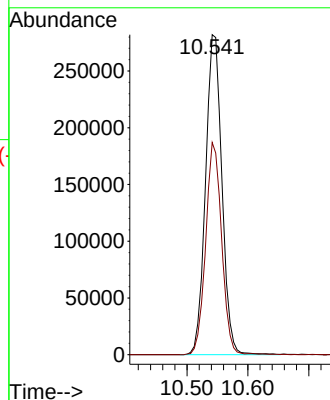
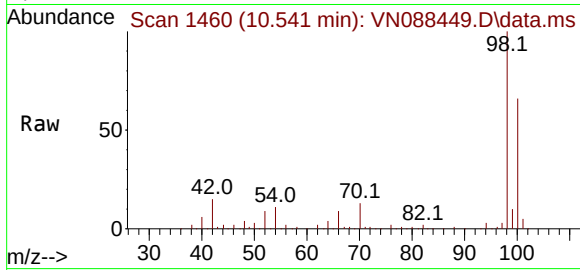




#50
Toluene-d8
Concen: 45.047 ug/l
RT: 10.541 min Scan# 1460
Delta R.T. -0.006 min
Lab File: VN088449.D
Acq: 04 Dec 2025 16:28

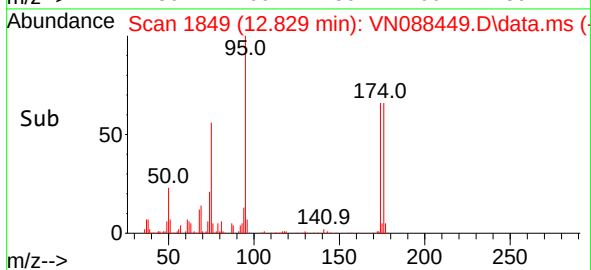
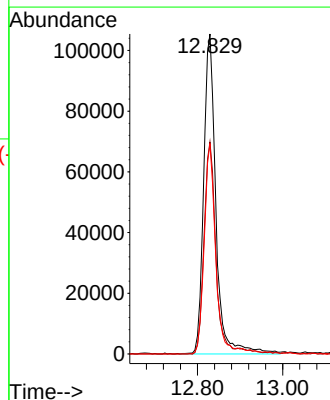
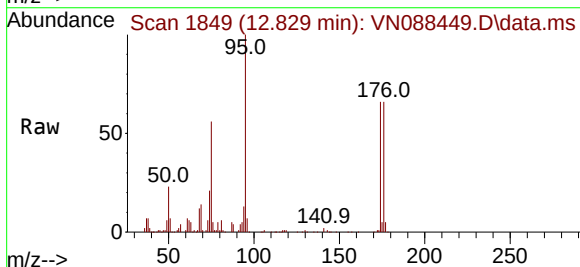
Instrument :
MSVOA_N
ClientSampleId :
TB01-120325

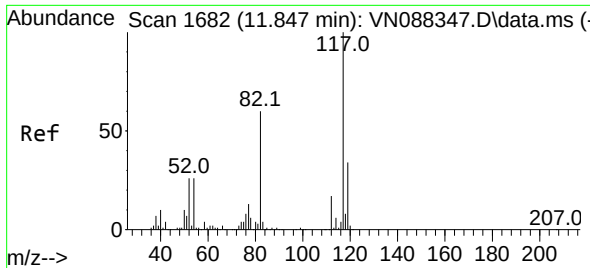
Tgt Ion: 98 Resp: 527723
Ion Ratio Lower Upper
98 100
100 64.6 52.2 78.2



#62
4-Bromofluorobenzene
Concen: 52.395 ug/l
RT: 12.829 min Scan# 1849
Delta R.T. 0.006 min
Lab File: VN088449.D
Acq: 04 Dec 2025 16:28

Tgt Ion: 95 Resp: 210610
Ion Ratio Lower Upper
95 100
174 62.5 0.0 139.0
176 62.3 0.0 132.4

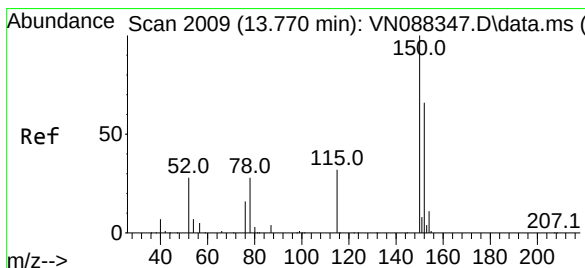
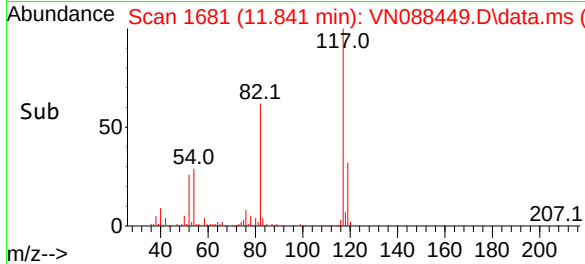
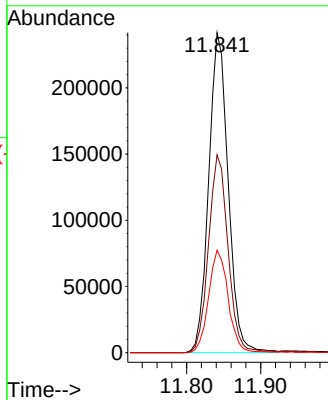
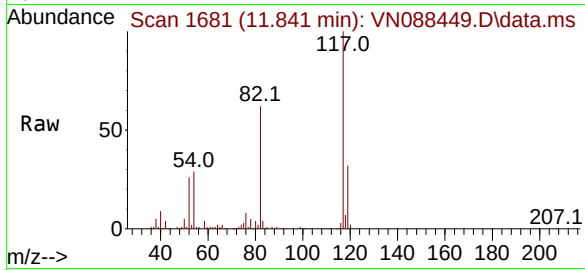




#63
Chlorobenzene-d5
Concen: 50.000 ug/l
RT: 11.841 min Scan# 1170
Delta R.T. -0.006 min
Lab File: VN088449.D
Acq: 04 Dec 2025 16:28

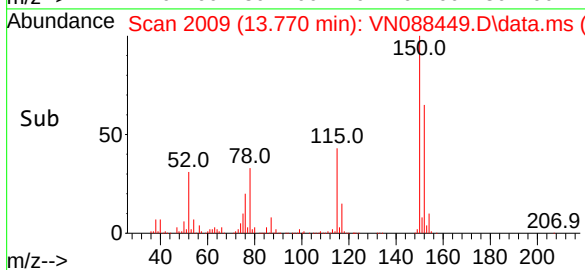
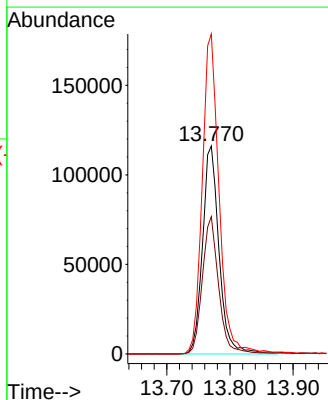
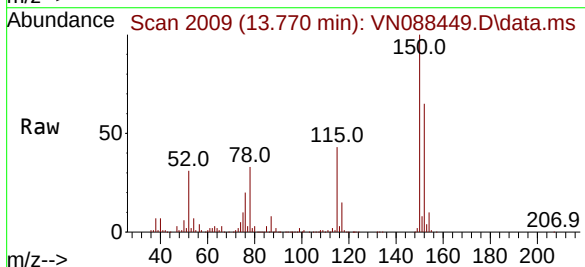
Instrument :
MSVOA_N
ClientSampleId :
TB01-120325

Tgt Ion	Ratio	Lower	Upper
117	100		
82	61.8	47.8	71.8
119	32.0	26.9	40.3



#72
1,4-Dichlorobenzene-d4
Concen: 50.000 ug/l
RT: 13.770 min Scan# 2009
Delta R.T. 0.000 min
Lab File: VN088449.D
Acq: 04 Dec 2025 16:28

Tgt Ion	Ratio	Lower	Upper
152	100		
115	65.0	33.0	98.9
150	156.8	0.0	349.0





CALIBRATION SUMMARY



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

VOLATILE ORGANICS INITIAL CALIBRATION DATA

Lab Name: Alliance Contract: JAC005
 Lab Code: ACE SDG No.: Q3757
 Instrument ID: MSVOA_N Calibration Date(s): 11/24/2025 11/24/2025
 Heated Purge: (Y/N) N Calibration Time(s): 12:38 14:35
 GC Column: RXI-624 ID: 0.25 (mm)

LAB FILE ID:		RRF001 = VN088344.D		RRF005 = VN088345.D		RRF020 = VN088346.D			
		RRF050 = VN088347.D		RRF100 = VN088348.D		RRF150 = VN088349.D			
COMPOUND	RRF001	RRF005	RRF020	RRF050	RRF100	RRF150	RRF	% RSD	
Vinyl Chloride	0.804	0.844	0.856	0.812	0.837	0.772	0.821	3.8	
1,1-Dichloroethene	0.722	0.620	0.601	0.568	0.595	0.554	0.610	9.8	
1,1-Dichloroethane	1.389	1.431	1.348	1.427	1.321	1.286	1.367	4.3	
cis-1,2-Dichloroethene	0.797	0.801	0.794	0.829	0.747	0.744	0.785	4.2	
1,1,1-Trichloroethane	1.247	1.219	1.192	1.158	1.166	1.104	1.181	4.3	
Benzene	1.658	1.555	1.609	1.559	1.559	1.477	1.570	3.9	
1,2-Dichloroethane	0.583	0.614	0.596	0.615	0.597	0.587	0.599	2.3	
Trichloroethene	0.382	0.385	0.383	0.357	0.371	0.348	0.371	4.1	
1,1,2-Trichloroethane	0.350	0.359	0.364	0.353	0.341	0.336	0.351	3	
Tetrachloroethene	0.426	0.388	0.413	0.377	0.407	0.367	0.396	5.7	
1,2-Dichloroethane-d4		0.954	0.916	0.999	0.882	0.960	0.942	4.7	
Dibromofluoromethane		0.361	0.350	0.350	0.334	0.337	0.346	3.2	
Toluene-d8		1.249	1.292	1.285	1.308	1.313	1.289	2	
4-Bromofluorobenzene		0.402	0.413	0.448	0.466	0.482	0.442	7.7	

* 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.

Method Path : Z:\voasrv\HPCHEM1\MSVOA_N\methods\
 Method File : 82N112425W.M
 Title : SW846 8260
 Last Update : Tue Nov 25 01:34:52 2025
 Response Via : Initial Calibration

Calibration Files

1 =VN088344.D 5 =VN088345.D 20 =VN088346.D 50 =VN088347.D 100 =VN088348.D 150 =VN088349.D

	Compound	1	5	20	50	100	150	Avg	%RSD

1) I	Pentafluorobenzene	-----ISTD-----							
2) T	Dichlorodifluo...	0.719	0.676	0.796	0.741	0.838	0.736	0.751	7.67
3) P	Chloromethane	0.911	0.800	0.805	0.801	0.759	0.764	0.807	6.79
4) C	Vinyl Chloride	0.804	0.844	0.856	0.812	0.837	0.772	0.821	3.78#
5) T	Bromomethane		0.407	0.392	0.396	0.387	0.392	0.395	1.85
6) T	Chloroethane	0.555	0.503	0.496	0.505	0.502	0.485	0.508	4.82
7) T	Trichlorofluor...	1.234	1.170	1.168	1.092	1.201	1.055	1.153	5.84
8) T	Diethyl Ether	0.495	0.451	0.395	0.444	0.401	0.411	0.433	8.73
9) T	1,1,2-Trichlor...	0.660	0.635	0.639	0.564	0.633	0.553	0.614	7.17
10) T	Methyl Iodide		0.498	0.584	0.662	0.660	0.636	0.608	11.40
11) T	Tert butyl alc...		0.151	0.149	0.161	0.142	0.148	0.150	4.64
12) CM	1,1-Dichloroet...	0.722	0.620	0.601	0.568	0.595	0.554	0.610	9.83#
13) T	Acrolein		0.146	0.132	0.151	0.143	0.147	0.144	5.00
14) T	Allyl chloride	1.107	1.143	1.117	1.149	1.132	1.135	1.131	1.38
15) T	Acrylonitrile	0.435	0.435	0.432	0.466	0.416	0.418	0.434	4.14
16) T	Acetone	0.412	0.355	0.322	0.335	0.305	0.314	0.341	11.49
17) T	Carbon Disulfide	2.061	2.043	2.043	1.983	1.996	1.833	1.993	4.22
18) T	Methyl Acetate	1.072	1.048	0.957	1.049	0.972	0.985	1.014	4.74
19) T	Methyl tert-bu...	2.126	2.239	2.245	2.504	2.265	2.333	2.285	5.53
20) T	Methylene Chlo...	0.797	0.752	0.704	0.756	0.686	0.678	0.729	6.44
21) T	trans-1,2-Dich...	0.752	0.668	0.676	0.686	0.647	0.626	0.676	6.39
22) T	Diisopropyl ether	2.267	2.415	2.439	2.580	2.359	2.424	2.414	4.26
23) T	Vinyl Acetate	1.711	1.879	1.997	2.137	1.943	2.002	1.945	7.35
24) P	1,1-Dichloroet...	1.389	1.431	1.348	1.427	1.321	1.286	1.367	4.27
25) T	2-Butanone	0.509	0.538	0.548	0.596	0.540	0.556	0.548	5.17
26) T	2,2-Dichloropr...	1.245	1.133	1.185	1.204	1.176	1.137	1.180	3.56
27) T	cis-1,2-Dichlo...	0.797	0.801	0.794	0.829	0.747	0.744	0.785	4.22
28) T	Bromochloromet...	0.724	0.722	0.684	0.735	0.624	0.597	0.681	8.47
29) T	Tetrahydrofuran	0.354	0.397	0.394	0.429	0.384	0.398	0.393	6.18
30) C	Chloroform	1.377	1.362	1.351	1.392	1.291	1.285	1.343	3.33#
31) T	Cyclohexane		1.450	1.264	1.133	1.237	1.109	1.239	10.92
32) T	1,1,1-Trichlor...	1.247	1.219	1.192	1.158	1.166	1.104	1.181	4.27
33) S	1,2-Dichloroet...		0.954	0.916	0.999	0.882	0.960	0.942	4.74
34) I	1,4-Difluorobenzene	-----ISTD-----							
35) S	Dibromofluorom...		0.361	0.350	0.350	0.334	0.337	0.346	3.17
36) T	1,1-Dichloropr...	0.607	0.542	0.531	0.491	0.547	0.490	0.535	8.09
37) T	Ethyl Acetate	0.735	0.678	0.724	0.720	0.681	0.681	0.703	3.69
38) T	Carbon Tetrach...	0.565	0.504	0.548	0.499	0.570	0.509	0.532	6.04
39) T	Methylcyclohexane	0.534	0.542	0.606	0.527	0.662	0.561	0.572	9.14
40) TM	Benzene	1.658	1.555	1.609	1.559	1.559	1.477	1.570	3.86
41) T	Methacrylonitrile	0.360	0.323	0.351	0.357	0.349	0.350	0.348	3.74
42) TM	1,2-Dichloroet...	0.583	0.614	0.596	0.615	0.597	0.587	0.599	2.27
43) T	Isopropyl Acetate	1.098	1.019	1.063	1.116	1.076	1.074	1.074	3.11
44) TM	Trichloroethene	0.382	0.385	0.383	0.357	0.371	0.348	0.371	4.13
45) C	1,2-Dichloropr...	0.426	0.402	0.402	0.406	0.396	0.381	0.402	3.62#
46) T	Dibromomethane	0.288	0.292	0.289	0.293	0.283	0.277	0.287	2.08
47) T	Bromodichlorom...	0.602	0.586	0.586	0.589	0.583	0.566	0.586	1.96
48) T	Methyl methacr...	0.480	0.433	0.491	0.513	0.511	0.513	0.490	6.35
49) T	1,4-Dioxane	0.005	0.006	0.007	0.007	0.006	0.006	0.006	8.55
50) S	Toluene-d8		1.249	1.292	1.285	1.308	1.313	1.289	1.96
51) T	4-Methyl-2-Pen...	0.540	0.616	0.661	0.673	0.658	0.651	0.633	7.79
52) CM	Toluene	0.829	0.885	0.940	0.913	0.960	0.911	0.906	5.05#
53) T	t-1,3-Dichloro...	0.567	0.553	0.598	0.631	0.631	0.625	0.601	5.72
54) T	cis-1,3-Dichlo...	0.590	0.618	0.628	0.662	0.654	0.640	0.632	4.16
55) T	1,1,2-Trichlor...	0.350	0.359	0.364	0.353	0.341	0.336	0.351	2.98
56) T	Ethyl methacry...	0.417	0.482	0.558	0.616	0.624	0.626	0.554	15.71

Method Path : Z:\voasrv\HPCHEM1\MSVOA_N\methods\
 Method File : 82N112425W.M

57)	T	1,3-Dichloropr...	0.607	0.638	0.647	0.644	0.638	0.622	0.633	2.42
58)	T	2-Chloroethyl ...	0.258	0.323	0.333	0.364	0.337	0.301	0.319	11.41
59)	T	2-Hexanone	0.359	0.301	0.406	0.436	0.442	0.450	0.399	14.66
60)	T	Dibromochlorom...	0.384	0.386	0.406	0.409	0.410	0.404	0.400	2.87
61)	T	1,2-Dibromoethane	0.304	0.367	0.364	0.369	0.367	0.359	0.355	7.16
62)	S	4-Bromofluorob...	0.402	0.413	0.448	0.466	0.482	0.442		7.67
63)	I	Chlorobenzene-d5	-----ISTD-----							
64)	T	Tetrachloroethene	0.426	0.388	0.413	0.377	0.407	0.367	0.396	5.73
65)	PM	Chlorobenzene	1.114	1.142	1.161	1.106	1.145	1.070	1.123	2.96
66)	T	1,1,1,2-Tetrac...	0.378	0.372	0.394	0.383	0.386	0.361	0.379	3.04
67)	C	Ethyl Benzene	1.881	1.860	2.014	1.938	2.133	1.952	1.963	5.06#
68)	T	m/p-Xylenes	0.658	0.690	0.769	0.733	0.788	0.728	0.728	6.64
69)	T	o-Xylene	0.670	0.619	0.722	0.696	0.749	0.708	0.694	6.51
70)	T	Styrene	0.884	1.025	1.163	1.193	1.284	1.210	1.127	12.94
71)	P	Bromoform	0.224	0.273	0.285	0.285	0.302	0.295	0.277	10.12
72)	I	1,4-Dichlorobenzen...	-----ISTD-----							
73)	T	Isopropylbenzene	3.339	3.528	3.944	3.624	4.040	3.706	3.697	7.07
74)	T	N-amyl acetate			1.173	1.045	1.338	1.480	1.259	15.09
75)	P	1,1,2,2-Tetrac...	1.377	1.218	1.231	1.127	1.135	1.082	1.195	8.87
76)	T	1,2,3-Trichlor...	1.036	1.199	1.268	1.202	1.177	1.145	1.171	6.63
77)	T	Bromobenzene	0.781	0.885	0.866	0.863	0.865	0.816	0.846	4.66
78)	T	n-propylbenzene	4.076	4.270	4.925	4.439	4.931	4.496	4.523	7.66
79)	T	2-Chlorotoluene	2.707	2.799	2.851	2.699	2.835	2.681	2.762	2.72
80)	T	1,3,5-Trimethy...	2.650	2.988	3.315	3.051	3.307	3.069	3.063	7.99
81)	T	trans-1,4-Dich...		0.355	0.364	0.430	0.452	0.462	0.413	12.09
82)	T	4-Chlorotoluene	2.093	2.830	2.935	2.854	2.944	2.803	2.743	11.80
83)	T	tert-Butylbenzene	2.509	2.468	2.773	2.509	2.816	2.549	2.604	5.78
84)	T	1,2,4-Trimethy...	2.572	2.827	3.310	3.112	3.344	3.105	3.045	9.74
85)	T	sec-Butylbenzene	3.457	3.614	3.971	3.609	4.122	3.674	3.741	6.73
86)	T	p-Isopropyltol...	2.714	2.808	3.330	3.047	3.435	3.073	3.068	9.18
87)	T	1,3-Dichlorobe...	1.730	1.573	1.720	1.606	1.676	1.574	1.647	4.34
88)	T	1,4-Dichlorobe...	1.747	1.825	1.786	1.681	1.702	1.608	1.725	4.52
89)	T	n-Butylbenzene	2.827	2.664	3.155	2.873	3.373	2.979	2.979	8.49
90)	T	Hexachloroethane	0.549	0.550	0.571	0.535	0.604	0.553	0.560	4.36
91)	T	1,2-Dichlorobe...	1.518	1.574	1.583	1.521	1.595	1.505	1.550	2.52
92)	T	1,2-Dibromo-3-...	0.284	0.284	0.284	0.288	0.300	0.300	0.290	2.80
93)	T	1,2,4-Trichlor...	0.998	0.864	0.901	0.889	0.955	0.917	0.921	5.26
94)	T	Hexachlorobuta...	0.362	0.316	0.349	0.297	0.354	0.307	0.331	8.32
95)	T	Naphthalene	2.946	2.864	3.163	3.292	3.483	3.395	3.190	7.74
96)	T	1,2,3-Trichlor...	0.942	0.850	0.888	0.875	0.930	0.891	0.896	3.83

(#) = Out of Range

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN112425\
 Data File : VN088344.D
 Acq On : 24 Nov 2025 12:38
 Operator : JC\MD
 Sample : VSTDICC001
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 VSTDICC001

Manual Integrations
 APPROVED

Reviewed By :John
 Carlone

Quant Time: Nov 25 01:09:22 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N112425W.M
 Quant Title : SW846 8260
 QLast Update : Tue Nov 25 01:08:44 2025
 Response via : Initial Calibration

11/25/2025
 Supervised By :Mahesh
 Dadoda

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

Internal Standards						
1) Pentafluorobenzene	8.206	168	286338	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.082	114	535294	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.841	117	458943	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.770	152	203454	50.000	ug/l	0.00

11/26/2025

System Monitoring Compounds

33) 1,2-Dichloroethane-d4	0.000	65	0d	0.000	ug/l	
Spiked Amount	50.000	Range	74 - 125	Recovery	=	0.000%#
35) Dibromofluoromethane	0.000	113	0d	0.000	ug/l	
Spiked Amount	50.000	Range	75 - 124	Recovery	=	0.000%#
50) Toluene-d8	0.000	98	0d	0.000	ug/l	
Spiked Amount	50.000	Range	86 - 113	Recovery	=	0.000%#
62) 4-Bromofluorobenzene	0.000	95	0d	0.000	ug/l	
Spiked Amount	50.000	Range	77 - 121	Recovery	=	0.000%#

Target Compounds

					Qvalue	
2) Dichlorodifluoromethane	2.142	85	4116	1.176	ug/l	88
3) Chloromethane	2.383	50	5215	1.282	ug/l	92
4) Vinyl Chloride	2.536	62	4602	1.055	ug/l	98
6) Chloroethane	3.148	64	3181	1.036	ug/l #	78
7) Trichlorofluoromethane	3.512	101	7064	1.049	ug/l #	82
8) Diethyl Ether	3.953	74	2832	1.101	ug/l	94
9) 1,1,2-Trichlorotrifluo...	4.359	101	3781	1.083	ug/l	91
12) 1,1-Dichloroethene	4.336	96	4136m	1.118	ug/l	
14) Allyl chloride	5.012	41	6342	0.957	ug/l	95
15) Acrylonitrile	5.712	53	12465	5.068	ug/l	98
16) Acetone	4.424	43	11802	4.716	ug/l	98
17) Carbon Disulfide	4.694	76	11805	1.034	ug/l #	85
18) Methyl Acetate	5.006	43	6138	1.129	ug/l	95
19) Methyl tert-butyl Ether	5.783	73	12174	0.876	ug/l	96
20) Methylene Chloride	5.265	84	4566	1.042	ug/l #	70
21) trans-1,2-Dichloroethene	5.783	96	4308	1.053	ug/l #	94
22) Diisopropyl ether	6.665	45	12982m	0.932	ug/l	
23) Vinyl Acetate	6.594	43	48988	3.972	ug/l	96
24) 1,1-Dichloroethane	6.547	63	7954	1.029	ug/l #	95
25) 2-Butanone	7.471	43	14574	4.283	ug/l #	82
26) 2,2-Dichloropropane	7.465	77	7127	1.026	ug/l #	77
27) cis-1,2-Dichloroethene	7.465	96	4562m	0.962	ug/l	
28) Bromochloromethane	7.788	49	4148	1.149	ug/l #	92
29) Tetrahydrofuran	7.835	42	10143	4.482	ug/l	96
30) Chloroform	7.947	83	7885	1.027	ug/l	97
32) 1,1,1-Trichloroethane	8.153	97	7143	1.055	ug/l #	46
36) 1,1-Dichloropropene	8.341	75	6499	1.203	ug/l #	80
37) Ethyl Acetate	7.553	43	7870	1.095	ug/l #	94
38) Carbon Tetrachloride	8.347	117	6045	1.085	ug/l	84
39) Methylcyclohexane	9.582	83	5715	0.846	ug/l	87
40) Benzene	8.582	78	17750	1.069	ug/l #	88
41) Methacrylonitrile	7.759	41	3857	1.037	ug/l #	83
42) 1,2-Dichloroethane	8.659	62	6237	1.039	ug/l	99
43) Isopropyl Acetate	8.677	43	11759	1.073	ug/l #	94
44) Trichloroethene	9.335	130	4088	1.057	ug/l	73
45) 1,2-Dichloropropane	9.600	63	4559	1.082	ug/l #	78

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN112425\
 Data File : VN088344.D
 Acq On : 24 Nov 2025 12:38
 Operator : JC\MD
 Sample : VSTDICC001
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 VSTDICC001

Manual Integrations
 APPROVED

Reviewed By :John
 Carlone

11/25/2025
 Supervised By :Mahesh
 Dadoda

Quant Time: Nov 25 01:09:22 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N112425W.M
 Quant Title : SW846 8260
 QLast Update : Tue Nov 25 01:08:44 2025
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
46) Dibromomethane	9.694	93	3080	1.022	ug/l	88
47) Bromodichloromethane	9.865	83	6445	1.068	ug/l #	89
48) Methyl methacrylate	9.665	41	5144	1.017	ug/l #	86
49) 1,4-Dioxane	9.688	88	1133	17.773	ug/l #	72
51) 4-Methyl-2-Pentanone	10.429	43	28924	4.454	ug/l	97
52) Toluene	10.612	92	8872	0.875	ug/l	97
53) t-1,3-Dichloropropene	10.812	75	6068	0.947	ug/l	89
54) cis-1,3-Dichloropropene	10.294	75	6312	0.928	ug/l	94
55) 1,1,2-Trichloroethane	11.000	97	3748	0.969	ug/l #	72
56) Ethyl methacrylate	10.859	69	4463	0.723	ug/l #	72
57) 1,3-Dichloropropane	11.147	76	6496	0.952	ug/l	95
58) 2-Chloroethyl Vinyl ether	10.141	63	13790	4.190	ug/l	92
59) 2-Hexanone	11.188	43	19197m	4.495	ug/l	
60) Dibromochloromethane	11.335	129	4116	0.936	ug/l	98
61) 1,2-Dibromoethane	11.447	107	3251	0.814	ug/l	96
64) Tetrachloroethene	11.076	164	3914	1.220	ug/l #	86
65) Chlorobenzene	11.871	112	10226	0.958	ug/l #	84
66) 1,1,1,2-Tetrachloroethane	11.941	131	3473	0.985	ug/l #	57
67) Ethyl Benzene	11.941	91	17267	0.922	ug/l	92
68) m/p-Xylenes	12.047	106	12077	1.724	ug/l	99
69) o-Xylene	12.382	106	6153	0.927	ug/l	91
70) Styrene	12.394	104	8118	0.747	ug/l	100
71) Bromoform	12.559	173	2054	0.782	ug/l #	82
73) Isopropylbenzene	12.676	105	13585	0.849	ug/l	99
75) 1,1,2,2-Tetrachloroethane	12.918	83	5605	1.121	ug/l #	83
76) 1,2,3-Trichloropropane	12.982	75	4214m	0.871	ug/l	
77) Bromobenzene	12.965	156	3176	0.907	ug/l	100
78) n-propylbenzene	13.017	91	16584	0.851	ug/l	98
79) 2-Chlorotoluene	13.106	91	11013	0.981	ug/l	100
80) 1,3,5-Trimethylbenzene	13.153	105	10783	0.811	ug/l	96
82) 4-Chlorotoluene	13.212	91	8516	0.700	ug/l	84
83) tert-Butylbenzene	13.417	119	10210	0.869	ug/l	95
84) 1,2,4-Trimethylbenzene	13.465	105	10464	0.786	ug/l	89
85) sec-Butylbenzene	13.594	105	14068	0.819	ug/l	99
86) p-Isopropyltoluene	13.706	119	11043	0.798	ug/l	97
87) 1,3-Dichlorobenzene	13.712	146	7040	1.019	ug/l	95
88) 1,4-Dichlorobenzene	13.782	146	7109m	0.955	ug/l	
89) n-Butylbenzene	14.035	91	11505	0.845	ug/l	93
90) Hexachloroethane	14.312	117	2233	0.841	ug/l	98
91) 1,2-Dichlorobenzene	14.082	146	6178	0.927	ug/l	97
92) 1,2-Dibromo-3-Chloropr...	14.694	75	1155	0.973	ug/l	93
93) 1,2,4-Trichlorobenzene	15.376	180	4059	1.011	ug/l	91
94) Hexachlorobutadiene	15.476	225	1474	1.002	ug/l	93
95) Naphthalene	15.623	128	11988	0.863	ug/l #	93
96) 1,2,3-Trichlorobenzene	15.823	180	3833	0.989	ug/l	100

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

```
Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN112425\  
Data File : VN088344.D  
Acq On    : 24 Nov 2025 12:38  
Operator  : JC\MD  
Sample    : VSTDIC001  
Misc      : 5.0mL/MSVOA_N/WATER  
ALS Vial  : 2    Sample Multiplier: 1
```

Instrument :
MSVOA_N
ClientSampleId :
VSTDICC001

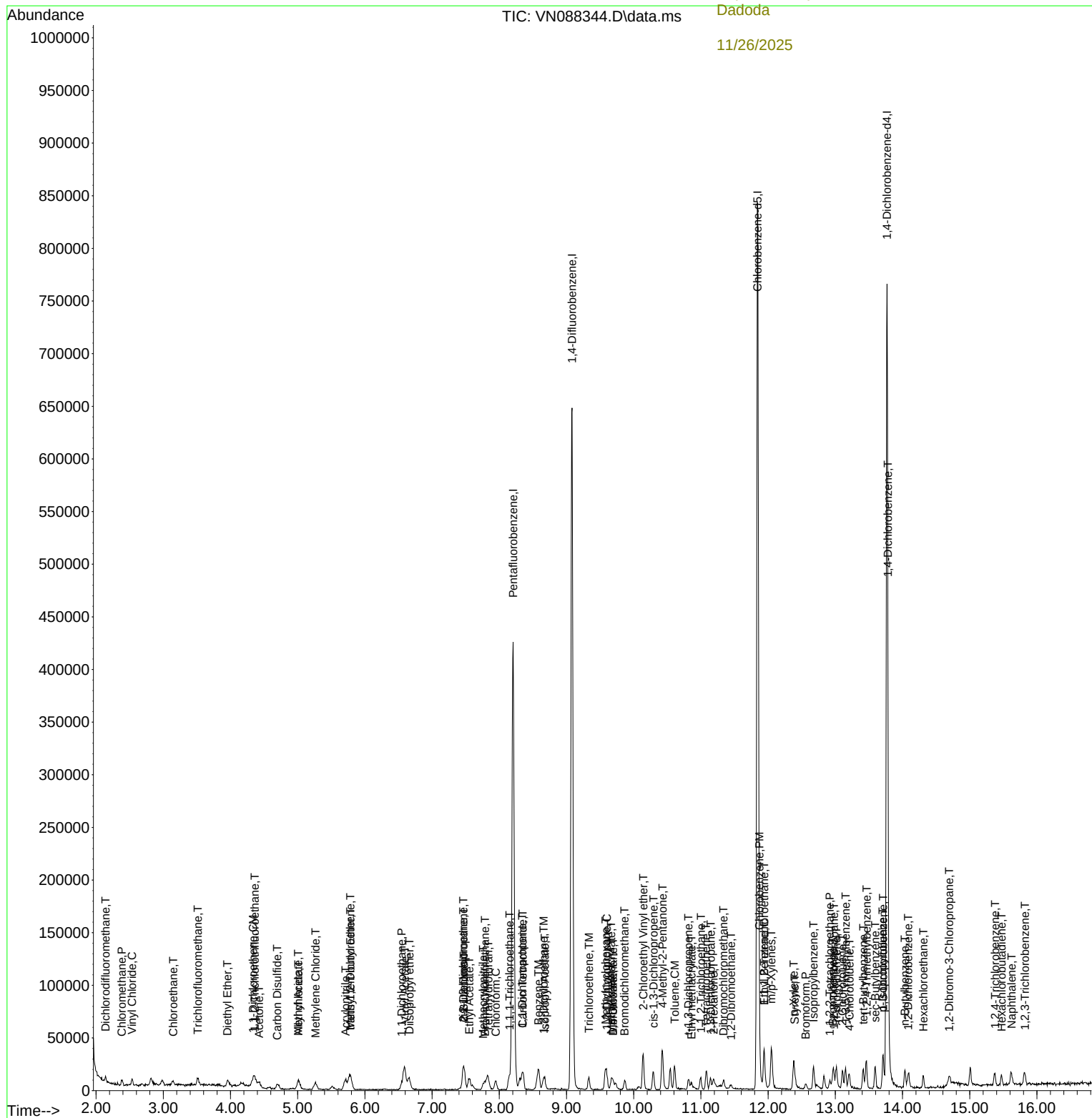
Manual Integrations APPROVED

Reviewed By :John
Carlone

11/25/2025
Supervised By :Mahesh
Dadoda

11/26/2025

Quant Time: Nov 25 01:09:22 2025
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N112425W.M
Quant Title : SW846 8260
QLast Update : Tue Nov 25 01:08:44 2025
Response via : Initial Calibration



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN112425\
 Data File : VN088345.D
 Acq On : 24 Nov 2025 13:11
 Operator : JC\MD
 Sample : VSTDICC005
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 VSTDICC005

Manual Integrations
 APPROVED

Reviewed By : John Carlone 11/25/2025
 Supervised By : Mahesh Dadoda 11/26/2025

Quant Time: Nov 25 01:10:10 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N112425W.M
 Quant Title : SW846 8260
 QLast Update : Tue Nov 25 01:08:44 2025
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.206	168	277017	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.083	114	520222	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.841	117	454370	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.771	152	210682	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.553	65	26416	5.272	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery =	10.540%#		
35) Dibromofluoromethane	8.153	113	18802	5.271	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery =	10.540%#		
50) Toluene-d8	10.547	98	64962	4.778	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery =	9.560%#		
62) 4-Bromofluorobenzene	12.829	95	20929	4.383	ug/l	0.00
Spiked Amount 50.000	Range 77 - 121		Recovery =	8.760%#		
Target Compounds					Qvalue	
2) Dichlorodifluoromethane	2.142	85	18728	5.530	ug/l	78
3) Chloromethane	2.383	50	22161	5.630	ug/l	96
4) Vinyl Chloride	2.536	62	23377	5.542	ug/l	95
5) Bromomethane	2.971	94	11265	4.586	ug/l	89
6) Chloroethane	3.142	64	13946	4.694	ug/l	90
7) Trichlorofluoromethane	3.512	101	32404	4.974	ug/l	99
8) Diethyl Ether	3.959	74	12500	5.022	ug/l	96
9) 1,1,2-Trichlorotrifluo...	4.365	101	17577	5.203	ug/l	97
10) Methyl Iodide	4.577	142	13782	4.252	ug/l #	86
11) Tert butyl alcohol	5.524	59	20975	23.946	ug/l	98
12) 1,1-Dichloroethene	4.342	96	17186	4.802	ug/l #	75
13) Acrolein	4.171	56	20281	21.167	ug/l	95
14) Allyl chloride	5.006	41	31674	4.939	ug/l	98
15) Acrylonitrile	5.706	53	60208	25.303	ug/l	99
16) Acetone	4.424	43	49209	20.327	ug/l	97
17) Carbon Disulfide	4.695	76	56594	5.124	ug/l	98
18) Methyl Acetate	5.012	43	29034	5.519	ug/l	99
19) Methyl tert-butyl Ether	5.783	73	62035	4.613	ug/l	96
20) Methylene Chloride	5.271	84	20834	4.915	ug/l #	80
21) trans-1,2-Dichloroethene	5.777	96	18495	4.671	ug/l	96
22) Diisopropyl ether	6.653	45	66886	4.963	ug/l	94
23) Vinyl Acetate	6.589	43	260201m	21.808	ug/l	
24) 1,1-Dichloroethane	6.547	63	39629	5.298	ug/l	98
25) 2-Butanone	7.465	43	74555	22.648	ug/l	97
26) 2,2-Dichloropropane	7.465	77	31387	4.669	ug/l	99
27) cis-1,2-Dichloroethene	7.471	96	22190	4.838	ug/l	96
28) Bromochloromethane	7.794	49	19995	5.723	ug/l	99
29) Tetrahydrofuran	7.824	42	54983	25.113	ug/l	100
30) Chloroform	7.953	83	37740	5.081	ug/l	99
31) Cyclohexane	8.241	56	40171	5.829	ug/l	96
32) 1,1,1-Trichloroethane	8.147	97	33772	5.154	ug/l #	85
36) 1,1-Dichloropropene	8.347	75	28193	5.368	ug/l	96
37) Ethyl Acetate	7.547	43	35286	5.054	ug/l	98
38) Carbon Tetrachloride	8.341	117	26194	4.839	ug/l	89
39) Methylcyclohexane	9.583	83	28179	4.292	ug/l	95
40) Benzene	8.588	78	80904	5.014	ug/l	91

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN112425\
 Data File : VN088345.D
 Acq On : 24 Nov 2025 13:11
 Operator : JC\MD
 Sample : VSTDICC005
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 3 Sample Multiplier: 1

Instrument :

MSVOA_N

ClientSampleId :

VSTDICC005

Manual Integrations

APPROVED

Reviewed By :John Carlone 11/25/2025

Supervised By :Mahesh Dadoda 11/26/2025

Quant Time: Nov 25 01:10:10 2025

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N112425W.M

Quant Title : SW846 8260

QLast Update : Tue Nov 25 01:08:44 2025

Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Methacrylonitrile	7.765	41	16819	4.652	ug/l	95
42) 1,2-Dichloroethane	8.653	62	31962	5.480	ug/l	98
43) Isopropyl Acetate	8.677	43	52990	4.976	ug/l	99
44) Trichloroethene	9.335	130	20005	5.321	ug/l	89
45) 1,2-Dichloropropane	9.600	63	20897	5.101	ug/l	96
46) Dibromomethane	9.688	93	15209	5.195	ug/l	98
47) Bromodichloromethane	9.865	83	30479	5.198	ug/l #	91
48) Methyl methacrylate	9.665	41	22513	4.579	ug/l	95
49) 1,4-Dioxane	9.688	88	6435	103.870	ug/l #	96
51) 4-Methyl-2-Pentanone	10.424	43	160328	25.401	ug/l	100
52) Toluene	10.606	92	46057	4.673	ug/l	99
53) t-1,3-Dichloropropene	10.812	75	28774	4.621	ug/l	98
54) cis-1,3-Dichloropropene	10.288	75	32170	4.869	ug/l	98
55) 1,1,2-Trichloroethane	10.994	97	18669	4.965	ug/l	96
56) Ethyl methacrylate	10.859	69	25089	4.184	ug/l	98
57) 1,3-Dichloropropane	11.141	76	33183	5.006	ug/l	96
58) 2-Chloroethyl Vinyl ether	10.141	63	83959	26.251	ug/l	98
59) 2-Hexanone	11.182	43	78248	18.852	ug/l	96
60) Dibromochloromethane	11.335	129	20078	4.697	ug/l	96
61) 1,2-Dibromoethane	11.447	107	19090	4.919	ug/l	95
64) Tetrachloroethene	11.077	164	17642	5.552	ug/l	92
65) Chlorobenzene	11.871	112	51903	4.912	ug/l	100
66) 1,1,1,2-Tetrachloroethane	11.935	131	16899	4.842	ug/l	95
67) Ethyl Benzene	11.941	91	84509	4.556	ug/l	97
68) m/p-Xylenes	12.047	106	62717	9.044	ug/l	96
69) o-Xylene	12.382	106	28111	4.278	ug/l	94
70) Styrene	12.394	104	46583	4.327	ug/l	99
71) Bromoform	12.565	173	12418	4.774	ug/l #	98
73) Isopropylbenzene	12.676	105	74324	4.486	ug/l	96
75) 1,1,2,2-Tetrachloroethane	12.918	83	25666	4.956	ug/l	99
76) 1,2,3-Trichloropropane	12.971	75	25268m	5.045	ug/l	
77) Bromobenzene	12.959	156	18654	5.143	ug/l	93
78) n-propylbenzene	13.018	91	89952	4.455	ug/l	97
79) 2-Chlorotoluene	13.106	91	58975	5.072	ug/l	99
80) 1,3,5-Trimethylbenzene	13.153	105	62949	4.574	ug/l	97
81) trans-1,4-Dichloro-2-b...	12.729	75	7483	4.106	ug/l #	78
82) 4-Chlorotoluene	13.200	91	59617	4.734	ug/l	97
83) tert-Butylbenzene	13.418	119	51988	4.275	ug/l	99
84) 1,2,4-Trimethylbenzene	13.465	105	59560	4.319	ug/l	99
85) sec-Butylbenzene	13.594	105	76137	4.278	ug/l	99
86) p-Isopropyltoluene	13.706	119	59155	4.127	ug/l	99
87) 1,3-Dichlorobenzene	13.712	146	33144	4.633	ug/l	97
88) 1,4-Dichlorobenzene	13.788	146	38453	4.987	ug/l	93
89) n-Butylbenzene	14.035	91	56132	3.979	ug/l	99
90) Hexachloroethane	14.312	117	11594	4.214	ug/l	98
91) 1,2-Dichlorobenzene	14.088	146	33169	4.804	ug/l	97
92) 1,2-Dibromo-3-Chloropr...	14.700	75	5979	4.862	ug/l	93
93) 1,2,4-Trichlorobenzene	15.370	180	18203	4.379	ug/l	99
94) Hexachlorobutadiene	15.476	225	6667	4.376	ug/l	95
95) Naphthalene	15.617	128	60329	4.195	ug/l	98
96) 1,2,3-Trichlorobenzene	15.817	180	17918	4.465	ug/l	96

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN112425\
Data File : VN088345.D
Acq On : 24 Nov 2025 13:11
Operator : JC\MD
Sample : VSTDICC005
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 3 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
VSTDICC005

Manual Integrations
APPROVED

Reviewed By :John Carlone 11/25/2025
Supervised By :Mahesh Dadoda 11/26/2025

Quant Time: Nov 25 01:10:10 2025
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N112425W.M
Quant Title : SW846 8260
QLast Update : Tue Nov 25 01:08:44 2025
Response via : Initial Calibration

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

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN112425\
 Data File : VN088345.D
 Acq On : 24 Nov 2025 13:11
 Operator : JC\MD
 Sample : VSTDIC005
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 3 Sample Multiplier: 1

Instrument :

MSVOA_N

ClientSampleId :

VSTDIC005

Manual Integrations

APPROVED

Reviewed By :John Carlone 11/25/2025

Supervised By :Mahesh Dadoda 11/26/2025

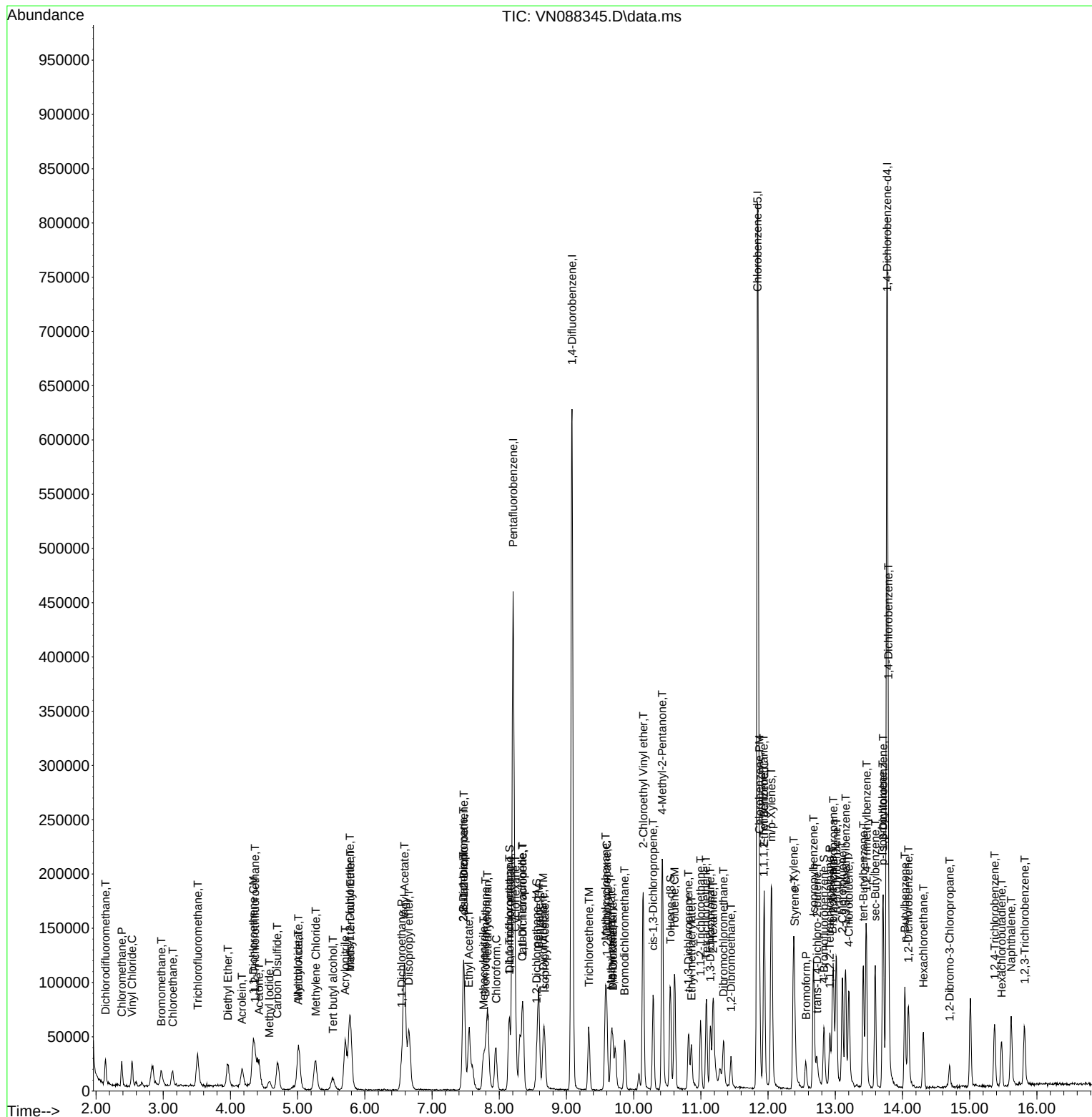
Quant Time: Nov 25 01:10:10 2025

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N112425W.M

Quant Title : SW846 8260

QLast Update : Tue Nov 25 01:08:44 2025

Response via : Initial Calibration



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN112425\
 Data File : VN088346.D
 Acq On : 24 Nov 2025 13:32
 Operator : JC\MD
 Sample : VSTDICC020
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 VSTDICC020

Manual Integrations
 APPROVED

Reviewed By : John Carlone 11/25/2025
 Supervised By : Mahesh Dadoda 11/26/2025

Quant Time: Nov 25 01:10:54 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N112425W.M
 Quant Title : SW846 8260
 QLast Update : Tue Nov 25 01:08:44 2025
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.206	168	289343	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.082	114	521967	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.847	117	446626	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.770	152	210470	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.565	65	106028	20.258	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery =	40.520%#		
35) Dibromofluoromethane	8.147	113	73055	20.412	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery =	40.820%#		
50) Toluene-d8	10.547	98	269703	19.772	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery =	39.540%#		
62) 4-Bromofluorobenzene	12.823	95	86319	18.015	ug/l	0.00
Spiked Amount 50.000	Range 77 - 121		Recovery =	36.020%#		
Target Compounds					Qvalue	
2) Dichlorodifluoromethane	2.142	85	92076	26.032	ug/l	100
3) Chloromethane	2.383	50	93182	22.664	ug/l	98
4) Vinyl Chloride	2.542	62	99078	22.487	ug/l	96
5) Bromomethane	2.983	94	45414	17.699	ug/l	100
6) Chloroethane	3.142	64	57451	18.515	ug/l	96
7) Trichlorofluoromethane	3.518	101	135233	19.875	ug/l	92
8) Diethyl Ether	3.965	74	45755	17.600	ug/l	98
9) 1,1,2-Trichlorotrifluo...	4.377	101	73956	20.959	ug/l	98
10) Methyl Iodide	4.583	142	67559	19.958	ug/l	92
11) Tert butyl alcohol	5.518	59	86464	94.506	ug/l	98
12) 1,1-Dichloroethene	4.341	96	69587	18.614	ug/l	85
13) Acrolein	4.177	56	76430	76.371	ug/l	97
14) Allyl chloride	5.012	41	129326	19.307	ug/l	97
15) Acrylonitrile	5.712	53	250269	100.698	ug/l	99
16) Acetone	4.424	43	186448	73.735	ug/l	96
17) Carbon Disulfide	4.706	76	236468	20.496	ug/l	96
18) Methyl Acetate	5.018	43	110778	20.161	ug/l	100
19) Methyl tert-butyl Ether	5.788	73	259864	18.502	ug/l	94
20) Methylene Chloride	5.265	84	81446	18.395	ug/l	97
21) trans-1,2-Dichloroethene	5.777	96	78223	18.913	ug/l	98
22) Diisopropyl ether	6.659	45	282262	20.050	ug/l	96
23) Vinyl Acetate	6.588	43	1155611m	92.730	ug/l	
24) 1,1-Dichloroethane	6.553	63	156012	19.969	ug/l	100
25) 2-Butanone	7.471	43	316859	92.154	ug/l	99
26) 2,2-Dichloropropane	7.471	77	137140	19.531	ug/l	99
27) cis-1,2-Dichloroethene	7.471	96	91877	19.178	ug/l	97
28) Bromochloromethane	7.794	49	79117	21.681	ug/l	100
29) Tetrahydrofuran	7.829	42	228189	99.785	ug/l	99
30) Chloroform	7.953	83	156345	20.153	ug/l	99
31) Cyclohexane	8.235	56	146307	20.327	ug/l	98
32) 1,1,1-Trichloroethane	8.153	97	137952	20.154	ug/l	98
36) 1,1-Dichloropropene	8.353	75	110824	21.032	ug/l	98
37) Ethyl Acetate	7.547	43	151153	21.576	ug/l	99
38) Carbon Tetrachloride	8.347	117	114374	21.060	ug/l	95
39) Methylcyclohexane	9.582	83	126502	19.203	ug/l	96
40) Benzene	8.588	78	335961	20.753	ug/l	98

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN112425\
 Data File : VN088346.D
 Acq On : 24 Nov 2025 13:32
 Operator : JC\MD
 Sample : VSTDICC020
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 4 Sample Multiplier: 1

Instrument :

MSVOA_N

ClientSampleId :

VSTDICC020

Manual Integrations

APPROVED

Reviewed By :John Carlone 11/25/2025

Supervised By :Mahesh Dadoda 11/26/2025

Quant Time: Nov 25 01:10:54 2025

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N112425W.M

Quant Title : SW846 8260

QLast Update : Tue Nov 25 01:08:44 2025

Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Methacrylonitrile	7.759	41	73287	20.203	ug/l	96
42) 1,2-Dichloroethane	8.653	62	124453	21.266	ug/l	98
43) Isopropyl Acetate	8.671	43	221861	20.765	ug/l	98
44) Trichloroethene	9.329	130	79876	21.175	ug/l	91
45) 1,2-Dichloropropane	9.600	63	83979	20.433	ug/l #	87
46) Dibromomethane	9.688	93	60253	20.513	ug/l	98
47) Bromodichloromethane	9.865	83	122415	20.808	ug/l	93
48) Methyl methacrylate	9.659	41	102560	20.792	ug/l	97
49) 1,4-Dioxane	9.676	88	28593	459.989	ug/l #	98
51) 4-Methyl-2-Pentanone	10.423	43	689576	108.887	ug/l	99
52) Toluene	10.612	92	196223	19.843	ug/l	99
53) t-1,3-Dichloropropene	10.818	75	124926	19.996	ug/l	99
54) cis-1,3-Dichloropropene	10.288	75	131162	19.787	ug/l	98
55) 1,1,2-Trichloroethane	10.994	97	75992	20.142	ug/l	96
56) Ethyl methacrylate	10.853	69	116597	19.380	ug/l	97
57) 1,3-Dichloropropane	11.141	76	135028	20.301	ug/l	99
58) 2-Chloroethyl Vinyl ether	10.141	63	347449	108.270	ug/l	98
59) 2-Hexanone	11.176	43	424244	101.867	ug/l	98
60) Dibromochloromethane	11.335	129	84716	19.753	ug/l	99
61) 1,2-Dibromoethane	11.447	107	76002	19.518	ug/l	99
64) Tetrachloroethene	11.076	164	73746	23.611	ug/l	96
65) Chlorobenzene	11.870	112	207406	19.968	ug/l	98
66) 1,1,1,2-Tetrachloroethane	11.935	131	70330	20.500	ug/l	98
67) Ethyl Benzene	11.941	91	359781	19.735	ug/l	96
68) m/p-Xylenes	12.047	106	274942	40.335	ug/l	99
69) o-Xylene	12.376	106	128919	19.957	ug/l	96
70) Styrene	12.388	104	207850	19.642	ug/l	99
71) Bromoform	12.559	173	50951	19.929	ug/l #	98
73) Isopropylbenzene	12.670	105	332053	20.064	ug/l	100
74) N-amyl acetate	12.682	43	98732m	20.946	ug/l	
75) 1,1,2,2-Tetrachloroethane	12.917	83	103664	20.037	ug/l	97
76) 1,2,3-Trichloropropane	12.970	75	106718m	21.331	ug/l	
77) Bromobenzene	12.959	156	72921	20.125	ug/l	98
78) n-propylbenzene	13.011	91	414648	20.557	ug/l	99
79) 2-Chlorotoluene	13.100	91	240017	20.663	ug/l	100
80) 1,3,5-Trimethylbenzene	13.147	105	279116	20.300	ug/l	98
81) trans-1,4-Dichloro-2-b...	12.717	75	30676	16.850	ug/l #	72
82) 4-Chlorotoluene	13.200	91	247131	19.645	ug/l	97
83) tert-Butylbenzene	13.417	119	233493	19.219	ug/l	99
84) 1,2,4-Trimethylbenzene	13.459	105	278702	20.229	ug/l	98
85) sec-Butylbenzene	13.594	105	334307	18.805	ug/l	99
86) p-Isopropyltoluene	13.706	119	280364	19.577	ug/l	99
87) 1,3-Dichlorobenzene	13.711	146	144838	20.267	ug/l	98
88) 1,4-Dichlorobenzene	13.788	146	150329	19.514	ug/l	99
89) n-Butylbenzene	14.035	91	265582	18.845	ug/l	99
90) Hexachloroethane	14.311	117	48067	17.489	ug/l	97
91) 1,2-Dichlorobenzene	14.088	146	133291	19.325	ug/l	98
92) 1,2-Dibromo-3-Chloropr...	14.694	75	23868	19.429	ug/l	99
93) 1,2,4-Trichlorobenzene	15.370	180	75854	18.265	ug/l	95
94) Hexachlorobutadiene	15.470	225	29359	19.288	ug/l	98
95) Naphthalene	15.611	128	266281	18.533	ug/l	99
96) 1,2,3-Trichlorobenzene	15.817	180	74795	18.658	ug/l	98

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN112425\
Data File : VN088346.D
Acq On : 24 Nov 2025 13:32
Operator : JC\MD
Sample : VSTDICC020
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 4 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
VSTDICC020

Manual Integrations
APPROVED

Reviewed By :John Carlone 11/25/2025
Supervised By :Mahesh Dadoda 11/26/2025

Quant Time: Nov 25 01:10:54 2025
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N112425W.M
Quant Title : SW846 8260
QLast Update : Tue Nov 25 01:08:44 2025
Response via : Initial Calibration

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

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

Instrument :
MSVOA_N
ClientSampleId :
VSTDICC020

Manual Integrations APPROVED

Reviewed By :John Carlone 11/25/2025
Supervised By :Mahesh Dadoda 11/26/2025



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN112425\
 Data File : VN088347.D
 Acq On : 24 Nov 2025 13:53
 Operator : JC\MD
 Sample : VSTDICCC050
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 VSTDICCC050

Manual Integrations
 APPROVED

Reviewed By : John Carlone 11/25/2025
 Supervised By : Mahesh Dadoda 11/26/2025

Quant Time: Nov 25 01:11:39 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N112425W.M
 Quant Title : SW846 8260
 QLast Update : Tue Nov 25 01:08:44 2025
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.206	168	254412	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.082	114	480905	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.847	117	425874	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.770	152	208615	50.000	ug/l	0.00

System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.559	65	254050	55.205	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	110.400%
35) Dibromofluoromethane	8.153	113	168216	51.013	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	102.020%
50) Toluene-d8	10.547	98	618062	49.180	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	98.360%
62) 4-Bromofluorobenzene	12.823	95	215443	48.802	ug/l	0.00
Spiked Amount	50.000	Range	77 - 121	Recovery	=	97.600%

Target Compounds					Qvalue	
2) Dichlorodifluoromethane	2.142	85	188518	60.616	ug/l	100
3) Chloromethane	2.383	50	203776	56.367	ug/l	100
4) Vinyl Chloride	2.536	62	206493	53.300	ug/l	100
5) Bromomethane	2.977	94	100800	44.679	ug/l	100
6) Chloroethane	3.142	64	128383	47.055	ug/l	100
7) Trichlorofluoromethane	3.512	101	277710	46.418	ug/l	100
8) Diethyl Ether	3.959	74	112943	49.408	ug/l	100
9) 1,1,2-Trichlorotrifluo...	4.365	101	143560	46.270	ug/l	100
10) Methyl Iodide	4.589	142	168409	56.580	ug/l	100
11) Tert butyl alcohol	5.512	59	205132	254.997	ug/l	100
12) 1,1-Dichloroethene	4.336	96	144499	43.959	ug/l	100
13) Acrolein	4.177	56	191880	218.057	ug/l	100
14) Allyl chloride	5.012	41	292223	49.615	ug/l	100
15) Acrylonitrile	5.706	53	593191	271.445	ug/l	100
16) Acetone	4.418	43	425884	191.549	ug/l	100
17) Carbon Disulfide	4.706	76	504570	49.738	ug/l	100
18) Methyl Acetate	5.018	43	266978	55.261	ug/l	100
19) Methyl tert-butyl Ether	5.783	73	637167	51.594	ug/l	100
20) Methylene Chloride	5.265	84	192304	49.396	ug/l	100
21) trans-1,2-Dichloroethene	5.771	96	174454	47.971	ug/l	100
22) Diisopropyl ether	6.659	45	656293	53.019	ug/l	100
23) Vinyl Acetate	6.588	43	2718764m	248.116	ug/l	
24) 1,1-Dichloroethane	6.559	63	362987	52.841	ug/l	100
25) 2-Butanone	7.465	43	757688	250.619	ug/l	100
26) 2,2-Dichloropropane	7.471	77	306308	49.612	ug/l	100
27) cis-1,2-Dichloroethene	7.471	96	210902	50.066	ug/l	100
28) Bromochloromethane	7.794	49	186871	58.241	ug/l	100
29) Tetrahydrofuran	7.824	42	545947	271.517	ug/l	100
30) Chloroform	7.947	83	354218	51.927	ug/l	100
31) Cyclohexane	8.235	56	288267	45.549	ug/l	100
32) 1,1,1-Trichloroethane	8.153	97	294555	48.942	ug/l	100
36) 1,1-Dichloropropene	8.353	75	236320	48.678	ug/l	100
37) Ethyl Acetate	7.547	43	346328	53.657	ug/l	100
38) Carbon Tetrachloride	8.347	117	240067	47.979	ug/l	100
39) Methylcyclohexane	9.582	83	253638	41.789	ug/l	100
40) Benzene	8.588	78	749637	50.260	ug/l	100

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN112425\
 Data File : VN088347.D
 Acq On : 24 Nov 2025 13:53
 Operator : JC\MD
 Sample : VSTDICCC050
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 VSTDICCC050

Manual Integrations
 APPROVED

Reviewed By : John Carlone 11/25/2025
 Supervised By : Mahesh Dadoda 11/26/2025

Quant Time: Nov 25 01:11:39 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N112425W.M
 Quant Title : SW846 8260
 QLast Update : Tue Nov 25 01:08:44 2025
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Methacrylonitrile	7.759	41	171571	51.334	ug/l	100
42) 1,2-Dichloroethane	8.653	62	295993	54.896	ug/l	100
43) Isopropyl Acetate	8.671	43	536716	54.522	ug/l	100
44) Trichloroethene	9.329	130	171828	49.440	ug/l	100
45) 1,2-Dichloropropane	9.600	63	195340	51.585	ug/l	100
46) Dibromomethane	9.688	93	140875	52.056	ug/l	100
47) Bromodichloromethane	9.871	83	283447	52.293	ug/l	100
48) Methyl methacrylate	9.659	41	246767	54.300	ug/l	100
49) 1,4-Dioxane	9.676	88	63666	1111.677	ug/l #	100
51) 4-Methyl-2-Pentanone	10.423	43	1618000	277.304	ug/l	100
52) Toluene	10.606	92	439167	48.202	ug/l	100
53) t-1,3-Dichloropropene	10.812	75	303537	52.732	ug/l	100
54) cis-1,3-Dichloropropene	10.294	75	318360	52.127	ug/l	100
55) 1,1,2-Trichloroethane	10.994	97	169966	48.898	ug/l	100
56) Ethyl methacrylate	10.853	69	296224	53.440	ug/l	100
57) 1,3-Dichloropropane	11.141	76	309692	50.536	ug/l	100
58) 2-Chloroethyl Vinyl ether	10.141	63	875006	295.946	ug/l	100
59) 2-Hexanone	11.176	43	1048582	273.278	ug/l	100
60) Dibromochloromethane	11.335	129	196756	49.794	ug/l	100
61) 1,2-Dibromoethane	11.447	107	177630	49.512	ug/l	100
64) Tetrachloroethene	11.082	164	160566	53.914	ug/l	100
65) Chlorobenzene	11.870	112	471088	47.564	ug/l	100
66) 1,1,1,2-Tetrachloroethane	11.941	131	162993	49.826	ug/l	100
67) Ethyl Benzene	11.941	91	825262	47.473	ug/l	100
68) m/p-Xylenes	12.047	106	624414	96.067	ug/l	100
69) o-Xylene	12.376	106	296307	48.105	ug/l	100
70) Styrene	12.388	104	508032	50.349	ug/l	100
71) Bromoform	12.559	173	121394	49.795	ug/l #	100
73) Isopropylbenzene	12.670	105	756073	46.091	ug/l	100
74) N-amyl acetate	12.500	43	218031	46.667	ug/l	100
75) 1,1,2,2-Tetrachloroethane	12.917	83	235147	45.856	ug/l	100
76) 1,2,3-Trichloropropane	12.970	75	250728m	50.561	ug/l	
77) Bromobenzene	12.959	156	179949	50.105	ug/l	100
78) n-propylbenzene	13.012	91	926041	46.319	ug/l	100
79) 2-Chlorotoluene	13.100	91	563004	48.901	ug/l	100
80) 1,3,5-Trimethylbenzene	13.153	105	636506	46.704	ug/l	100
81) trans-1,4-Dichloro-2-b...	12.717	75	89717	49.717	ug/l #	100
82) 4-Chlorotoluene	13.200	91	595484	47.756	ug/l	100
83) tert-Butylbenzene	13.412	119	523419	43.467	ug/l	100
84) 1,2,4-Trimethylbenzene	13.459	105	649135	47.534	ug/l	100
85) sec-Butylbenzene	13.594	105	752927	42.729	ug/l	100
86) p-Isopropyltoluene	13.706	119	635661	44.782	ug/l	100
87) 1,3-Dichlorobenzene	13.711	146	334974	47.290	ug/l	100
88) 1,4-Dichlorobenzene	13.788	146	350784	45.940	ug/l	100
89) n-Butylbenzene	14.035	91	599399	42.911	ug/l	100
90) Hexachloroethane	14.306	117	111560	40.952	ug/l	100
91) 1,2-Dichlorobenzene	14.082	146	317314	46.414	ug/l	100
92) 1,2-Dibromo-3-Chloropr...	14.694	75	60137	49.387	ug/l	100
93) 1,2,4-Trichlorobenzene	15.364	180	185506	45.066	ug/l	100
94) Hexachlorobutadiene	15.470	225	61955	41.064	ug/l	100
95) Naphthalene	15.611	128	686671	48.217	ug/l	100
96) 1,2,3-Trichlorobenzene	15.811	180	182475	45.925	ug/l	100

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN112425\
Data File : VN088347.D
Acq On : 24 Nov 2025 13:53
Operator : JC\MD
Sample : VSTDICCC050
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 5 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
VSTDICCC050

Manual Integrations
APPROVED

Reviewed By :John Carlone 11/25/2025
Supervised By :Mahesh Dadoda 11/26/2025

Quant Time: Nov 25 01:11:39 2025
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N112425W.M
Quant Title : SW846 8260
QLast Update : Tue Nov 25 01:08:44 2025
Response via : Initial Calibration

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

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

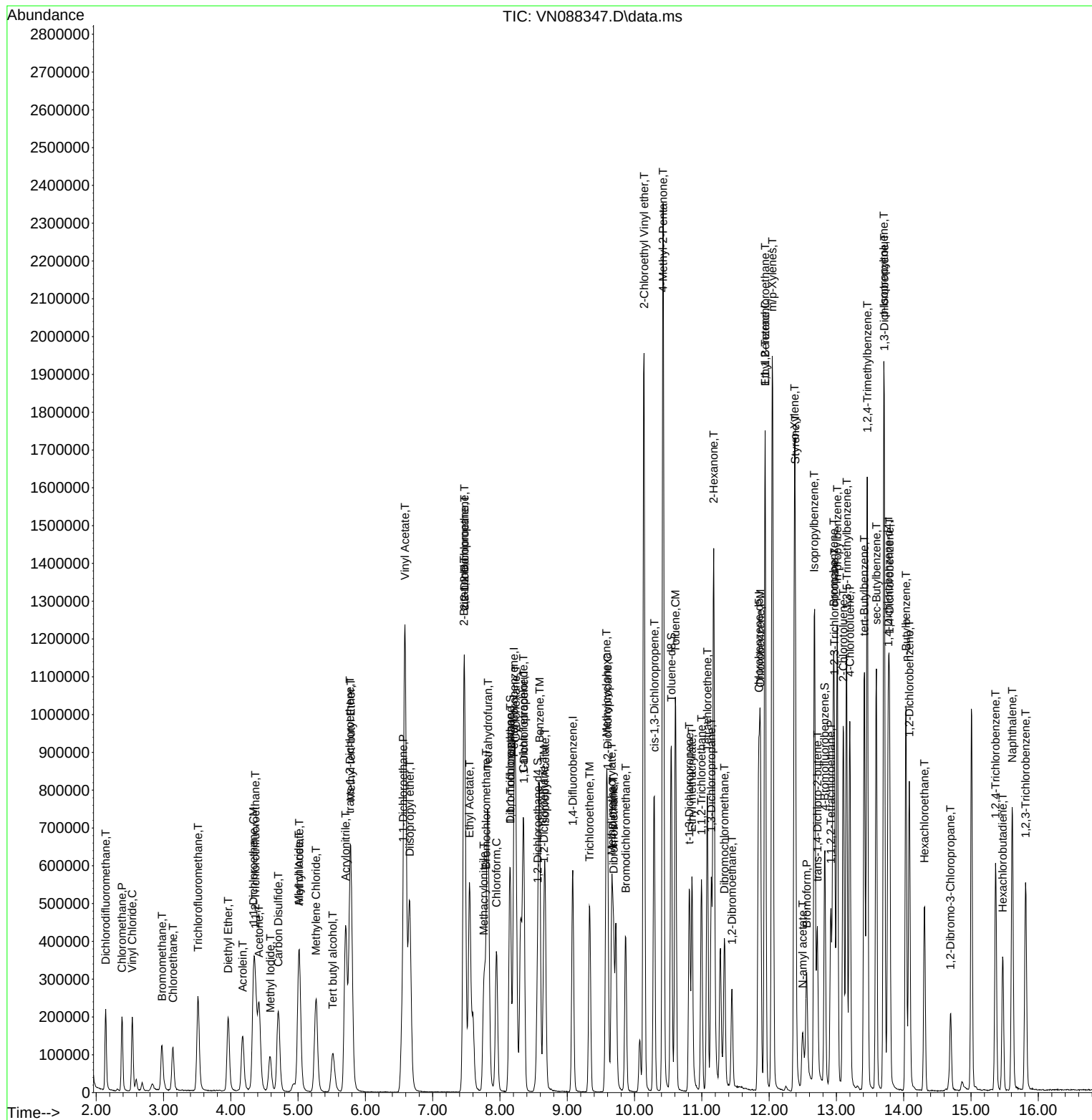
Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN112425\
Data File : VN088347.D
Acq On : 24 Nov 2025 13:53
Operator : JC\MD
Sample : VSTDICCC050
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 5 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
VSTDICCC050

Manual Integrations APPROVED

Reviewed By :John Carlone 11/25/2025
Supervised By :Mahesh Dadoda 11/26/2025

Quant Time: Nov 25 01:11:39 2025
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N112425W.M
Quant Title : SW846 8260
QLast Update : Tue Nov 25 01:08:44 2025
Response via : Initial Calibration



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN112425\
 Data File : VN088348.D
 Acq On : 24 Nov 2025 14:14
 Operator : JC\MD
 Sample : VSTDICC100
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 VSTDICC100

Quant Time: Nov 25 01:12:24 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N112425W.M
 Quant Title : SW846 8260
 QLast Update : Tue Nov 25 01:08:44 2025
 Response via : Initial Calibration

Manual Integrations APPROVED

Reviewed By : John Carlone 11/25/2025
 Supervised By : Mahesh Dadoda 11/26/2025

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

Internal Standards						
1) Pentafluorobenzene	8.206	168	285895	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.083	114	501695	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.847	117	439316	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.770	152	219473	50.000	ug/l	0.00

System Monitoring Compounds

33) 1,2-Dichloroethane-d4	8.559	65	504045	97.467	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	= 194.940%#	
35) Dibromofluoromethane	8.147	113	335348	97.484	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 194.960%#	
50) Toluene-d8	10.547	98	1312124	100.081	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	= 200.160%#	
62) 4-Bromofluorobenzene	12.829	95	467895	101.595	ug/l	0.00
Spiked Amount	50.000	Range	77 - 121	Recovery	= 203.200%#	

Target Compounds

						Qvalue
2) Dichlorodifluoromethane	2.142	85	479358	137.160	ug/l	100
3) Chloromethane	2.383	50	434116	106.858	ug/l	98
4) Vinyl Chloride	2.536	62	478320	109.869	ug/l	98
5) Bromomethane	2.965	94	221311	87.292	ug/l	99
6) Chloroethane	3.130	64	287091	93.637	ug/l	96
7) Trichlorofluoromethane	3.506	101	686944	102.176	ug/l	99
8) Diethyl Ether	3.959	74	229525	89.352	ug/l	99
9) 1,1,2-Trichlorotrifluo...	4.365	101	361718	103.745	ug/l	100
10) Methyl Iodide	4.577	142	377533	112.872	ug/l	97
11) Tert butyl alcohol	5.524	59	406566	449.741	ug/l	100
12) 1,1-Dichloroethene	4.336	96	340010	92.046	ug/l	96
13) Acrolein	4.177	56	409205	413.820	ug/l	99
14) Allyl chloride	5.012	41	647046m	97.762	ug/l	
15) Acrylonitrile	5.706	53	1189524	484.386	ug/l	99
16) Acetone	4.418	43	872591	349.245	ug/l	98
17) Carbon Disulfide	4.706	76	1141513	100.133	ug/l	99
18) Methyl Acetate	5.018	43	555796	102.374	ug/l	100
19) Methyl tert-butyl Ether	5.783	73	1295144	93.325	ug/l	100
20) Methylene Chloride	5.265	84	392179	89.644	ug/l	99
21) trans-1,2-Dichloroethene	5.777	96	369984	90.533	ug/l	98
22) Diisopropyl ether	6.659	45	1349029	96.981	ug/l	97
23) Vinyl Acetate	6.589	43	5554188m	451.061	ug/l	
24) 1,1-Dichloroethane	6.553	63	755476	97.866	ug/l	98
25) 2-Butanone	7.471	43	1542855	454.129	ug/l	97
26) 2,2-Dichloropropane	7.477	77	672523	96.933	ug/l	100
27) cis-1,2-Dichloroethene	7.471	96	427126	90.230	ug/l	98
28) Bromochloromethane	7.800	49	356867	98.975	ug/l	99
29) Tetrahydrofuran	7.824	42	1096740	485.379	ug/l	99
30) Chloroform	7.947	83	738399	96.326	ug/l	100
31) Cyclohexane	8.241	56	707339	99.459	ug/l	99
32) 1,1,1-Trichloroethane	8.147	97	666887	98.605	ug/l	98
36) 1,1-Dichloropropene	8.353	75	549319	108.461	ug/l	99
37) Ethyl Acetate	7.547	43	682857	101.412	ug/l	99
38) Carbon Tetrachloride	8.341	117	571861	109.553	ug/l	97
39) Methylcyclohexane	9.577	83	663987	104.863	ug/l	96
40) Benzene	8.588	78	1564367	100.538	ug/l	99

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN112425\
 Data File : VN088348.D
 Acq On : 24 Nov 2025 14:14
 Operator : JC\MD
 Sample : VSTDICC100
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 VSTDICC100

Manual Integrations
 APPROVED

Reviewed By : John Carlone 11/25/2025
 Supervised By : Mahesh Dadoda 11/26/2025

Quant Time: Nov 25 01:12:24 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N112425W.M
 Quant Title : SW846 8260
 QLast Update : Tue Nov 25 01:08:44 2025
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Methacrylonitrile	7.759	41	349832	100.333	ug/l	98
42) 1,2-Dichloroethane	8.653	62	599011	106.490	ug/l	100
43) Isopropyl Acetate	8.671	43	1079559	105.122	ug/l	99
44) Trichloroethene	9.335	130	372568	102.756	ug/l	98
45) 1,2-Dichloropropane	9.600	63	396928	100.477	ug/l	96
46) Dibromomethane	9.688	93	284059	100.615	ug/l	99
47) Bromodichloromethane	9.871	83	585372	103.520	ug/l	100
48) Methyl methacrylate	9.659	41	512503	108.100	ug/l	99
49) 1,4-Dioxane	9.682	88	124110	2077.291	ug/l #	98
51) 4-Methyl-2-Pentanone	10.424	43	3302590	542.564	ug/l	100
52) Toluene	10.606	92	962754	101.291	ug/l	98
53) t-1,3-Dichloropropene	10.818	75	633621	105.515	ug/l	99
54) cis-1,3-Dichloropropene	10.294	75	656506	103.040	ug/l	97
55) 1,1,2-Trichloroethane	10.994	97	342175	94.361	ug/l	97
56) Ethyl methacrylate	10.853	69	625749	108.210	ug/l	99
57) 1,3-Dichloropropane	11.141	76	640348	100.162	ug/l	99
58) 2-Chloroethyl Vinyl ether	10.141	63	1692209	548.624	ug/l	100
59) 2-Hexanone	11.177	43	2217550	553.982	ug/l	99
60) Dibromochloromethane	11.335	129	410889	99.677	ug/l	98
61) 1,2-Dibromoethane	11.447	107	368152	98.366	ug/l	100
64) Tetrachloroethene	11.082	164	357404	116.335	ug/l	95
65) Chlorobenzene	11.871	112	1006420	98.505	ug/l	99
66) 1,1,1,2-Tetrachloroethane	11.941	131	338748	100.385	ug/l	99
67) Ethyl Benzene	11.941	91	1873691	104.486	ug/l	99
68) m/p-Xylenes	12.047	106	1384206	206.446	ug/l	97
69) o-Xylene	12.376	106	657970	103.552	ug/l	99
70) Styrene	12.388	104	1127962	108.366	ug/l	99
71) Bromoform	12.559	173	265738	105.669	ug/l #	98
73) Isopropylbenzene	12.671	105	1773457	102.764	ug/l	100
74) N-amyl acetate	12.488	43	587472	119.520	ug/l	95
75) 1,1,2,2-Tetrachloroethane	12.918	83	498007	92.311	ug/l	100
76) 1,2,3-Trichloropropane	12.971	75	516579m	99.018	ug/l	
77) Bromobenzene	12.959	156	379553	100.455	ug/l	99
78) n-propylbenzene	13.012	91	2164268	102.898	ug/l	99
79) 2-Chlorotoluene	13.106	91	1244218	102.723	ug/l	99
80) 1,3,5-Trimethylbenzene	13.153	105	1451677	101.248	ug/l	99
81) trans-1,4-Dichloro-2-b...	12.718	75	198381	104.496	ug/l #	88
82) 4-Chlorotoluene	13.200	91	1292288	98.511	ug/l	100
83) tert-Butylbenzene	13.418	119	1236180	97.579	ug/l	99
84) 1,2,4-Trimethylbenzene	13.459	105	1467849	102.168	ug/l	98
85) sec-Butylbenzene	13.594	105	1809286	97.599	ug/l	100
86) p-Isopropyltoluene	13.706	119	1507703	100.961	ug/l	99
87) 1,3-Dichlorobenzene	13.712	146	735647	98.718	ug/l	99
88) 1,4-Dichlorobenzene	13.788	146	746880	92.975	ug/l	99
89) n-Butylbenzene	14.035	91	1480732	100.761	ug/l	99
90) Hexachloroethane	14.312	117	265251	92.554	ug/l	99
91) 1,2-Dichlorobenzene	14.082	146	700211	97.354	ug/l	99
92) 1,2-Dibromo-3-Chloropr...	14.700	75	131734	102.833	ug/l	99
93) 1,2,4-Trichlorobenzene	15.364	180	419347	96.834	ug/l	97
94) Hexachlorobutadiene	15.476	225	155408	97.909	ug/l	97
95) Naphthalene	15.612	128	1528952	102.050	ug/l	99
96) 1,2,3-Trichlorobenzene	15.812	180	408194	97.651	ug/l	98

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN112425\
Data File : VN088348.D
Acq On : 24 Nov 2025 14:14
Operator : JC\MD
Sample : VSTDICC100
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
VSTDICC100

Manual Integrations
APPROVED

Reviewed By :John Carlone 11/25/2025
Supervised By :Mahesh Dadoda 11/26/2025

Quant Time: Nov 25 01:12:24 2025
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N112425W.M
Quant Title : SW846 8260
QLast Update : Tue Nov 25 01:08:44 2025
Response via : Initial Calibration

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

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

```
Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN112425\  
Data File : VN088348.D  
Acq On    : 24 Nov 2025  14:14  
Operator  : JC\MD  
Sample    : VSTDICC100  
Misc      : 5.0mL/MSVOA_N/WATER  
ALS Vial  : 6    Sample Multiplier: 1
```

Instrument :

MSVOA N

ClientSampleId :

VSTDICC100

Quant Time: Nov 25 01:12:24 2025

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N112425W.M

Quant Title : SW846 8260

QLast Update : Tue Nov 25 01:08:44 2025

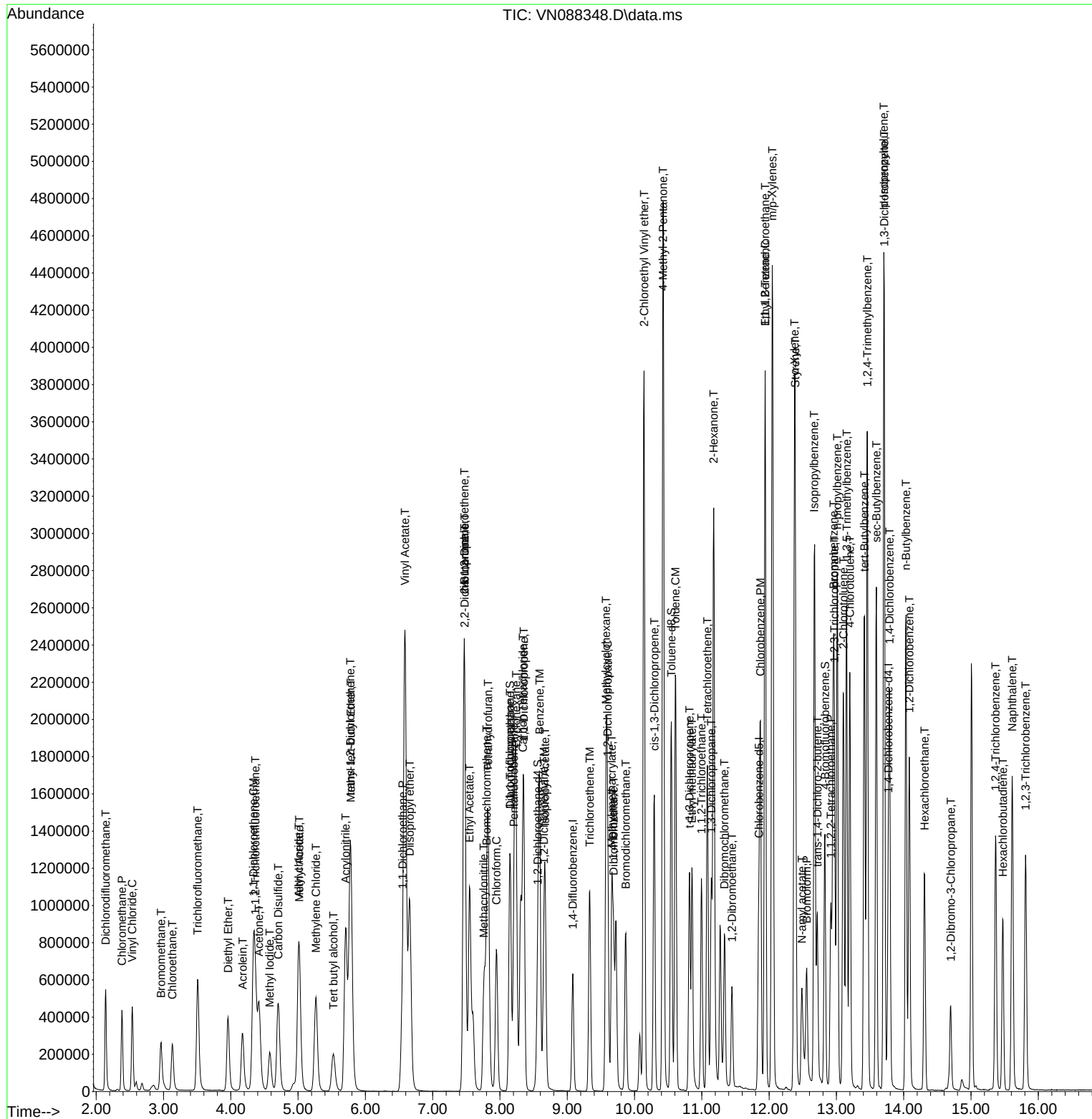
Response via : Initial Calibration

Manual Integrations

APPROVED

Reviewed By :John Carlone 11/25/2025

Supervised By :Mahesh Dadoda 11/26/2025



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN112425\
 Data File : VN088349.D
 Acq On : 24 Nov 2025 14:35
 Operator : JC\MD
 Sample : VSTDICC150
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 VSTDICC150

Manual Integrations
 APPROVED

Reviewed By : John Carlone 11/25/2025
 Supervised By : Mahesh Dadoda 11/26/2025

Quant Time: Nov 25 01:13:19 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N112425W.M
 Quant Title : SW846 8260
 QLast Update : Tue Nov 25 01:08:44 2025
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.206	168	278091	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.083	114	508326	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.847	117	457447	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.770	152	226673	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.559	65	800654	159.167	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery = 318.340%#			
35) Dibromofluoromethane	8.147	113	514160	147.513	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery = 295.020%#			
50) Toluene-d8	10.547	98	2002008	150.709	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery = 301.420%#			
62) 4-Bromofluorobenzene	12.829	95	734676	157.441	ug/l	0.00
Spiked Amount 50.000	Range 77 - 121		Recovery = 314.880%#			
Target Compounds						
						Qvalue
2) Dichlorodifluoromethane	2.142	85	614298	180.703	ug/l	97
3) Chloromethane	2.383	50	637070	161.216	ug/l	98
4) Vinyl Chloride	2.536	62	643786	152.025	ug/l	99
5) Bromomethane	2.948	94	327351	132.741	ug/l	99
6) Chloroethane	3.124	64	404469	135.623	ug/l	99
7) Trichlorofluoromethane	3.506	101	880178	134.591	ug/l	97
8) Diethyl Ether	3.959	74	342692	137.150	ug/l	100
9) 1,1,2-Trichlorotrifluo...	4.365	101	461517	136.083	ug/l	99
10) Methyl Iodide	4.583	142	530186	162.959	ug/l	96
11) Tert butyl alcohol	5.524	59	615624	700.111	ug/l	99
12) 1,1-Dichloroethene	4.336	96	461833	128.534	ug/l	95
13) Acrolein	4.177	56	614784	639.165	ug/l	98
14) Allyl chloride	5.012	41	947252m	147.136	ug/l	
15) Acrylonitrile	5.706	53	1745154	730.587	ug/l	100
16) Acetone	4.418	43	1310595	539.272	ug/l	100
17) Carbon Disulfide	4.700	76	1529256	137.911	ug/l	99
18) Methyl Acetate	5.018	43	822075	155.669	ug/l	100
19) Methyl tert-butyl Ether	5.783	73	1946194	144.173	ug/l	97
20) Methylene Chloride	5.259	84	565609	132.915	ug/l	99
21) trans-1,2-Dichloroethene	5.771	96	522186	131.362	ug/l	100
22) Diisopropyl ether	6.659	45	2022262	149.459	ug/l	98
23) Vinyl Acetate	6.589	43	8352858m	697.381	ug/l	
24) 1,1-Dichloroethane	6.553	63	1073175	142.923	ug/l	98
25) 2-Butanone	7.471	43	2317712	701.347	ug/l	96
26) 2,2-Dichloropropane	7.477	77	948461	140.541	ug/l	99
27) cis-1,2-Dichloroethene	7.471	96	621100	134.889	ug/l	99
28) Bromochloromethane	7.794	49	498418	142.113	ug/l	100
29) Tetrahydrofuran	7.824	42	1661991	756.181	ug/l	99
30) Chloroform	7.947	83	1072381	143.821	ug/l	95
31) Cyclohexane	8.236	56	925517	133.789	ug/l	98
32) 1,1,1-Trichloroethane	8.153	97	920717	139.956	ug/l	96
36) 1,1-Dichloropropene	8.353	75	747080	145.584	ug/l	100
37) Ethyl Acetate	7.547	43	1038194	152.172	ug/l	100
38) Carbon Tetrachloride	8.341	117	776182	146.756	ug/l	96
39) Methylcyclohexane	9.583	83	856176	133.452	ug/l	96
40) Benzene	8.588	78	2252306	142.862	ug/l	98

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN112425\
 Data File : VN088349.D
 Acq On : 24 Nov 2025 14:35
 Operator : JC\MD
 Sample : VSTDICC150
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 VSTDICC150

Manual Integrations
 APPROVED

Reviewed By : John Carlone 11/25/2025
 Supervised By : Mahesh Dadoda 11/26/2025

Quant Time: Nov 25 01:13:19 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N112425W.M
 Quant Title : SW846 8260
 QLast Update : Tue Nov 25 01:08:44 2025
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Methacrylonitrile	7.759	41	533051	150.887	ug/l	98
42) 1,2-Dichloroethane	8.653	62	895667	157.152	ug/l	99
43) Isopropyl Acetate	8.671	43	1638317	157.450	ug/l	99
44) Trichloroethene	9.335	130	529981	144.265	ug/l	98
45) 1,2-Dichloropropane	9.600	63	581336	145.238	ug/l	95
46) Dibromomethane	9.688	93	422737	147.782	ug/l	100
47) Bromodichloromethane	9.871	83	863631	150.736	ug/l	100
48) Methyl methacrylate	9.665	41	781838	162.758	ug/l	99
49) 1,4-Dioxane	9.683	88	191377	3161.388	ug/l #	97
51) 4-Methyl-2-Pentanone	10.424	43	4963362	804.767	ug/l	99
52) Toluene	10.612	92	1389323	144.264	ug/l	99
53) t-1,3-Dichloropropene	10.818	75	953623	156.732	ug/l	99
54) cis-1,3-Dichloropropene	10.294	75	975877	151.168	ug/l	98
55) 1,1,2-Trichloroethane	10.994	97	513073	139.644	ug/l	99
56) Ethyl methacrylate	10.853	69	954630	162.929	ug/l	99
57) 1,3-Dichloropropane	11.141	76	948252	146.389	ug/l	99
58) 2-Chloroethyl Vinyl ether	10.141	63	2298552	735.483	ug/l	100
59) 2-Hexanone	11.177	43	3427514	845.082	ug/l	99
60) Dibromochloromethane	11.335	129	615359	147.332	ug/l	98
61) 1,2-Dibromoethane	11.447	107	548222	144.568	ug/l	99
64) Tetrachloroethene	11.082	164	503663	157.444	ug/l	98
65) Chlorobenzene	11.871	112	1467895	137.978	ug/l	100
66) 1,1,1,2-Tetrachloroethane	11.941	131	494738	140.800	ug/l	98
67) Ethyl Benzene	11.941	91	2679329	143.490	ug/l	99
68) m/p-Xylenes	12.047	106	1999338	286.370	ug/l	98
69) o-Xylene	12.376	106	971419	146.823	ug/l	99
70) Styrene	12.388	104	1660474	153.203	ug/l	99
71) Bromoform	12.559	173	404703	154.550	ug/l #	99
73) Isopropylbenzene	12.671	105	2520465	141.411	ug/l	99
74) N-amyl acetate	12.482	43	1006304	198.228	ug/l #	95
75) 1,1,2,2-Tetrachloroethane	12.918	83	735899	132.074	ug/l	100
76) 1,2,3-Trichloropropane	12.971	75	778772m	144.533	ug/l	
77) Bromobenzene	12.959	156	554789	142.170	ug/l	97
78) n-propylbenzene	13.012	91	3057578	140.752	ug/l	100
79) 2-Chlorotoluene	13.100	91	1823307	145.751	ug/l	100
80) 1,3,5-Trimethylbenzene	13.153	105	2086957	140.933	ug/l	99
81) trans-1,4-Dichloro-2-b...	12.718	75	314465	160.381	ug/l #	91
82) 4-Chlorotoluene	13.200	91	1905793	140.664	ug/l	99
83) tert-Butylbenzene	13.418	119	1733611	132.498	ug/l	99
84) 1,2,4-Trimethylbenzene	13.459	105	2111285	142.286	ug/l	99
85) sec-Butylbenzene	13.594	105	2498053	130.473	ug/l	100
86) p-Isopropyltoluene	13.706	119	2089505	135.476	ug/l	99
87) 1,3-Dichlorobenzene	13.712	146	1070574	139.099	ug/l	100
88) 1,4-Dichlorobenzene	13.788	146	1093471	131.796	ug/l	99
89) n-Butylbenzene	14.035	91	2025634	133.462	ug/l	99
90) Hexachloroethane	14.306	117	376335	127.143	ug/l	99
91) 1,2-Dichlorobenzene	14.082	146	1023417	137.771	ug/l	100
92) 1,2-Dibromo-3-Chloropr...	14.700	75	204222	154.355	ug/l	100
93) 1,2,4-Trichlorobenzene	15.365	180	623560	139.416	ug/l	99
94) Hexachlorobutadiene	15.470	225	208495	127.182	ug/l	97
95) Naphthalene	15.612	128	2308447	149.183	ug/l	99
96) 1,2,3-Trichlorobenzene	15.812	180	605584	140.271	ug/l	98

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN112425\
Data File : VN088349.D
Acq On : 24 Nov 2025 14:35
Operator : JC\MD
Sample : VSTDICC150
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
VSTDICC150

Manual Integrations
APPROVED

Reviewed By :John Carlone 11/25/2025
Supervised By :Mahesh Dadoda 11/26/2025

Quant Time: Nov 25 01:13:19 2025
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N112425W.M
Quant Title : SW846 8260
QLast Update : Tue Nov 25 01:08:44 2025
Response via : Initial Calibration

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

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

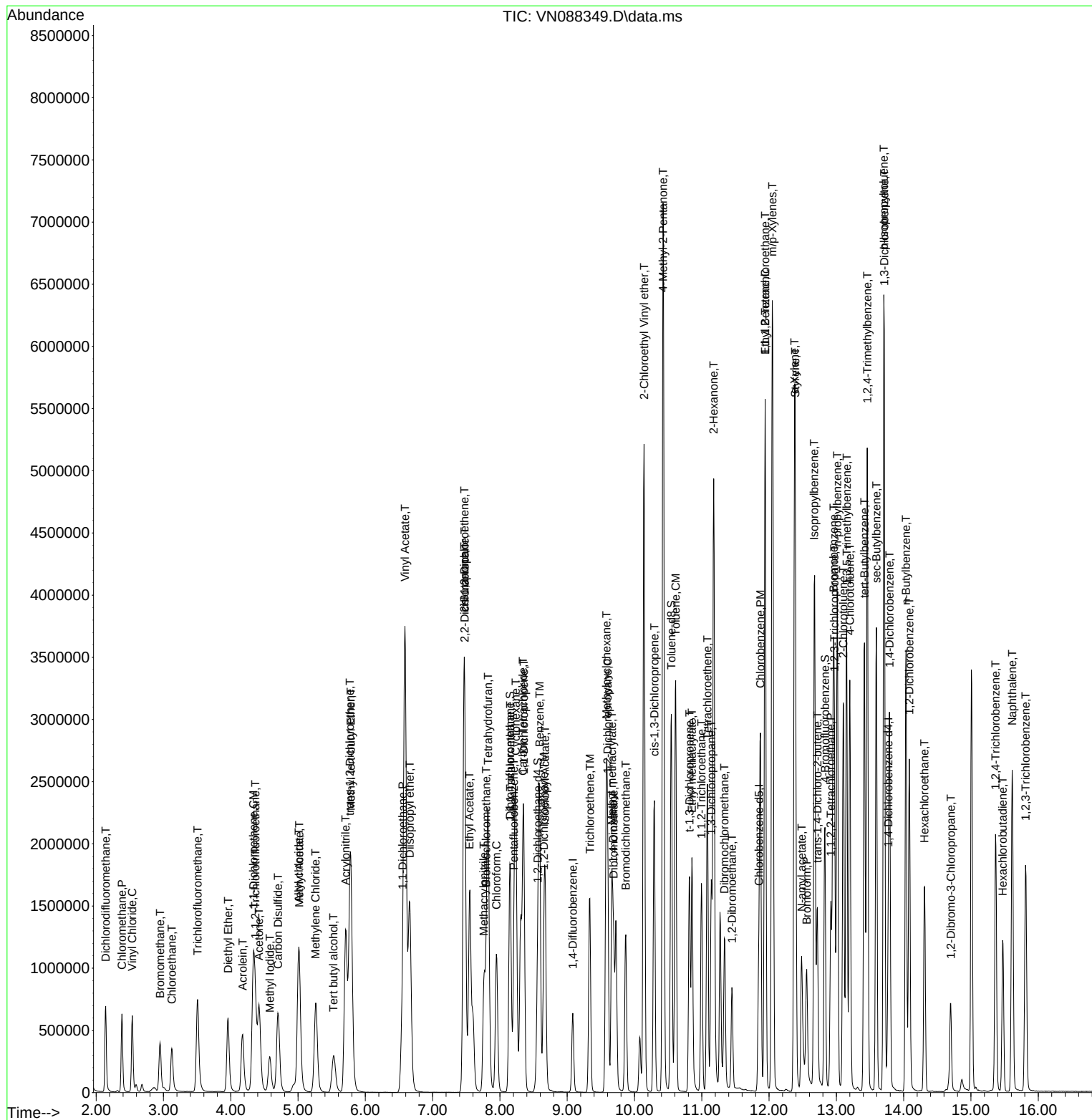

```
Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN112425\  
Data File : VN088349.D  
Acq On    : 24 Nov 2025  14:35  
Operator  : JC\MD  
Sample    : VSTDICC150  
Misc      : 5.0mL/MSVOA_N/WATER  
ALS Vial  : 7    Sample Multiplier: 1
```

Instrument :
MSVOA_N
ClientSampleId :
VSTDICC150

Manual Integrations APPROVED

Reviewed By :John Carlone 11/25/2025
Supervised By :Mahesh Dadoda 11/26/2025

Quant Time: Nov 25 01:13:19 2025
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N112425W.M
Quant Title : SW846 8260
QLast Update : Tue Nov 25 01:08:44 2025
Response via : Initial Calibration



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN112425\
 Data File : VN088351.D
 Acq On : 24 Nov 2025 15:18
 Operator : JC\MD
 Sample : VSTDICV050
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 ICVVN112425

Manual Integrations
 APPROVED

Reviewed By : John Carlone 11/25/2025
 Supervised By : Mahesh Dadoda 11/26/2025

Quant Time: Nov 25 01:39:30 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N112425W.M
 Quant Title : SW846 8260
 QLast Update : Tue Nov 25 01:34:52 2025
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.206	168	264693	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.082	114	504031	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.841	117	443578	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.770	152	211823	50.000	ug/l	0.00

System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.559	65	234844	47.098	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	94.200%
35) Dibromofluoromethane	8.153	113	152840	43.757	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	87.520%
50) Toluene-d8	10.547	98	574351	44.193	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	88.380%
62) 4-Bromofluorobenzene	12.829	95	199878	44.823	ug/l	0.00
Spiked Amount	50.000	Range	77 - 121	Recovery	=	89.640%

Target Compounds					Qvalue	
2) Dichlorodifluoromethane	2.142	85	204306	51.389	ug/l	96
3) Chloromethane	2.383	50	224506	52.578	ug/l	98
4) Vinyl Chloride	2.536	62	216512	49.842	ug/l	99
5) Bromomethane	2.977	94	113722	54.393	ug/l	95
6) Chloroethane	3.142	64	138310	51.450	ug/l	96
7) Trichlorofluoromethane	3.512	101	304206	49.826	ug/l	96
8) Diethyl Ether	3.965	74	122332	53.384	ug/l	99
9) 1,1,2-Trichlorotrifluo...	4.365	101	153067	47.094	ug/l	96
10) Methyl Iodide	4.589	142	173671	53.976	ug/l	95
11) Tert butyl alcohol	5.518	59	213569	268.272	ug/l	99
12) 1,1-Dichloroethene	4.342	96	158096	48.957	ug/l	94
13) Acrolein	4.177	56	203235	266.656	ug/l	99
14) Allyl chloride	5.018	41	321979	53.793	ug/l	99
15) Acrylonitrile	5.706	53	637644	277.614	ug/l	99
16) Acetone	4.418	43	450459	249.797	ug/l	100
17) Carbon Disulfide	4.706	76	544067	51.558	ug/l	100
18) Methyl Acetate	5.018	43	290918	54.196	ug/l	99
19) Methyl tert-butyl Ether	5.783	73	694497	57.401	ug/l	95
20) Methylene Chloride	5.265	84	208952	54.158	ug/l	98
21) trans-1,2-Dichloroethene	5.777	96	191102	53.421	ug/l	99
22) Diisopropyl ether	6.653	45	719883	56.335	ug/l	97
23) Vinyl Acetate	6.588	43	2915087m	283.141	ug/l	
24) 1,1-Dichloroethane	6.559	63	394512	54.516	ug/l	99
25) 2-Butanone	7.471	43	807706	278.613	ug/l	97
26) 2,2-Dichloropropane	7.471	77	337031	53.957	ug/l	99
27) cis-1,2-Dichloroethene	7.471	96	223944	53.866	ug/l	99
28) Bromochloromethane	7.794	49	167723	46.526	ug/l	100
29) Tetrahydrofuran	7.824	42	581218	279.515	ug/l	99
30) Chloroform	7.947	83	383675	53.957	ug/l	95
31) Cyclohexane	8.241	56	316159	48.211	ug/l	94
32) 1,1,1-Trichloroethane	8.147	97	319795	51.150	ug/l	100
36) 1,1-Dichloropropene	8.353	75	253183	46.967	ug/l	98
37) Ethyl Acetate	7.547	43	364546	51.431	ug/l	99
38) Carbon Tetrachloride	8.341	117	262078	48.837	ug/l	98
39) Methylcyclohexane	9.577	83	276272	47.913	ug/l	98
40) Benzene	8.588	78	803742	50.800	ug/l	100

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN112425\
 Data File : VN088351.D
 Acq On : 24 Nov 2025 15:18
 Operator : JC\MD
 Sample : VSTDICV050
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 ICVVN112425

Manual Integrations
 APPROVED

Reviewed By : John Carlone 11/25/2025
 Supervised By : Mahesh Dadoda 11/26/2025

Quant Time: Nov 25 01:39:30 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N112425W.M
 Quant Title : SW846 8260
 QLast Update : Tue Nov 25 01:34:52 2025
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Methacrylonitrile	7.765	41	189164	53.883	ug/l	98
42) 1,2-Dichloroethane	8.653	62	318757	52.806	ug/l	99
43) Isopropyl Acetate	8.671	43	575762	53.165	ug/l	100
44) Trichloroethene	9.335	130	184844	49.444	ug/l	97
45) 1,2-Dichloropropane	9.600	63	211177	52.095	ug/l	96
46) Dibromomethane	9.688	93	152417	52.686	ug/l	100
47) Bromodichloromethane	9.871	83	307856	52.155	ug/l	100
48) Methyl methacrylate	9.665	41	269474	54.535	ug/l	99
49) 1,4-Dioxane	9.682	88	64188	1021.452	ug/l #	99
51) 4-Methyl-2-Pentanone	10.424	43	1745522	273.447	ug/l	100
52) Toluene	10.612	92	475044	51.998	ug/l	100
53) t-1,3-Dichloropropene	10.812	75	323712	53.428	ug/l	97
54) cis-1,3-Dichloropropene	10.294	75	344582	54.081	ug/l	98
55) 1,1,2-Trichloroethane	10.994	97	182138	51.530	ug/l	98
56) Ethyl methacrylate	10.853	69	323702	57.977	ug/l	98
57) 1,3-Dichloropropane	11.141	76	338609	53.102	ug/l	98
58) 2-Chloroethyl Vinyl ether	10.141	63	851755	264.613	ug/l	100
59) 2-Hexanone	11.176	43	1118254	278.086	ug/l	99
60) Dibromochloromethane	11.341	129	212150	52.650	ug/l	98
61) 1,2-Dibromoethane	11.447	107	191658	53.546	ug/l	99
64) Tetrachloroethene	11.082	164	164101	46.666	ug/l	96
65) Chlorobenzene	11.871	112	512644	51.452	ug/l	100
66) 1,1,1,2-Tetrachloroethane	11.941	131	174200	51.838	ug/l	98
67) Ethyl Benzene	11.941	91	899198	51.635	ug/l	100
68) m/p-Xylenes	12.047	106	679183	105.191	ug/l	99
69) o-Xylene	12.376	106	322403	52.376	ug/l	100
70) Styrene	12.388	104	550033	55.031	ug/l	100
71) Bromoform	12.559	173	132448	53.811	ug/l #	100
73) Isopropylbenzene	12.670	105	827973	52.866	ug/l	100
74) N-amyl acetate	12.500	43	221082m	41.449	ug/l	
75) 1,1,2,2-Tetrachloroethane	12.917	83	253948	50.155	ug/l	100
76) 1,2,3-Trichloropropane	12.970	75	269266m	54.274	ug/l	
77) Bromobenzene	12.959	156	191381	53.406	ug/l	98
78) n-propylbenzene	13.012	91	1004887	52.446	ug/l	100
79) 2-Chlorotoluene	13.106	91	599091	51.202	ug/l	99
80) 1,3,5-Trimethylbenzene	13.153	105	687007	52.936	ug/l	99
81) trans-1,4-Dichloro-2-b...	12.717	75	96389	55.117	ug/l #	95
82) 4-Chlorotoluene	13.200	91	628637	54.093	ug/l	99
83) tert-Butylbenzene	13.417	119	569932	51.660	ug/l	99
84) 1,2,4-Trimethylbenzene	13.459	105	703272	54.519	ug/l	98
85) sec-Butylbenzene	13.594	105	812817	51.285	ug/l	100
86) p-Isopropyltoluene	13.706	119	678617	52.216	ug/l	98
87) 1,3-Dichlorobenzene	13.712	146	357734	51.282	ug/l	99
88) 1,4-Dichlorobenzene	13.788	146	372290	50.949	ug/l	98
89) n-Butylbenzene	14.035	91	644062	51.040	ug/l	99
90) Hexachloroethane	14.306	117	118158	49.768	ug/l	97
91) 1,2-Dichlorobenzene	14.082	146	344212	52.435	ug/l	99
92) 1,2-Dibromo-3-Chloropr...	14.700	75	62186	50.621	ug/l	98
93) 1,2,4-Trichlorobenzene	15.370	180	200095	51.301	ug/l	99
94) Hexachlorobutadiene	15.476	225	64430	45.969	ug/l	97
95) Naphthalene	15.611	128	719827	53.258	ug/l	99
96) 1,2,3-Trichlorobenzene	15.811	180	188353	49.620	ug/l	97

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN112425\
Data File : VN088351.D
Acq On : 24 Nov 2025 15:18
Operator : JC\MD
Sample : VSTDICV050
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
ICVVN112425

Manual Integrations
APPROVED

Reviewed By :John Carlone 11/25/2025
Supervised By :Mahesh Dadoda 11/26/2025

Quant Time: Nov 25 01:39:30 2025
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N112425W.M
Quant Title : SW846 8260
QLast Update : Tue Nov 25 01:34:52 2025
Response via : Initial Calibration

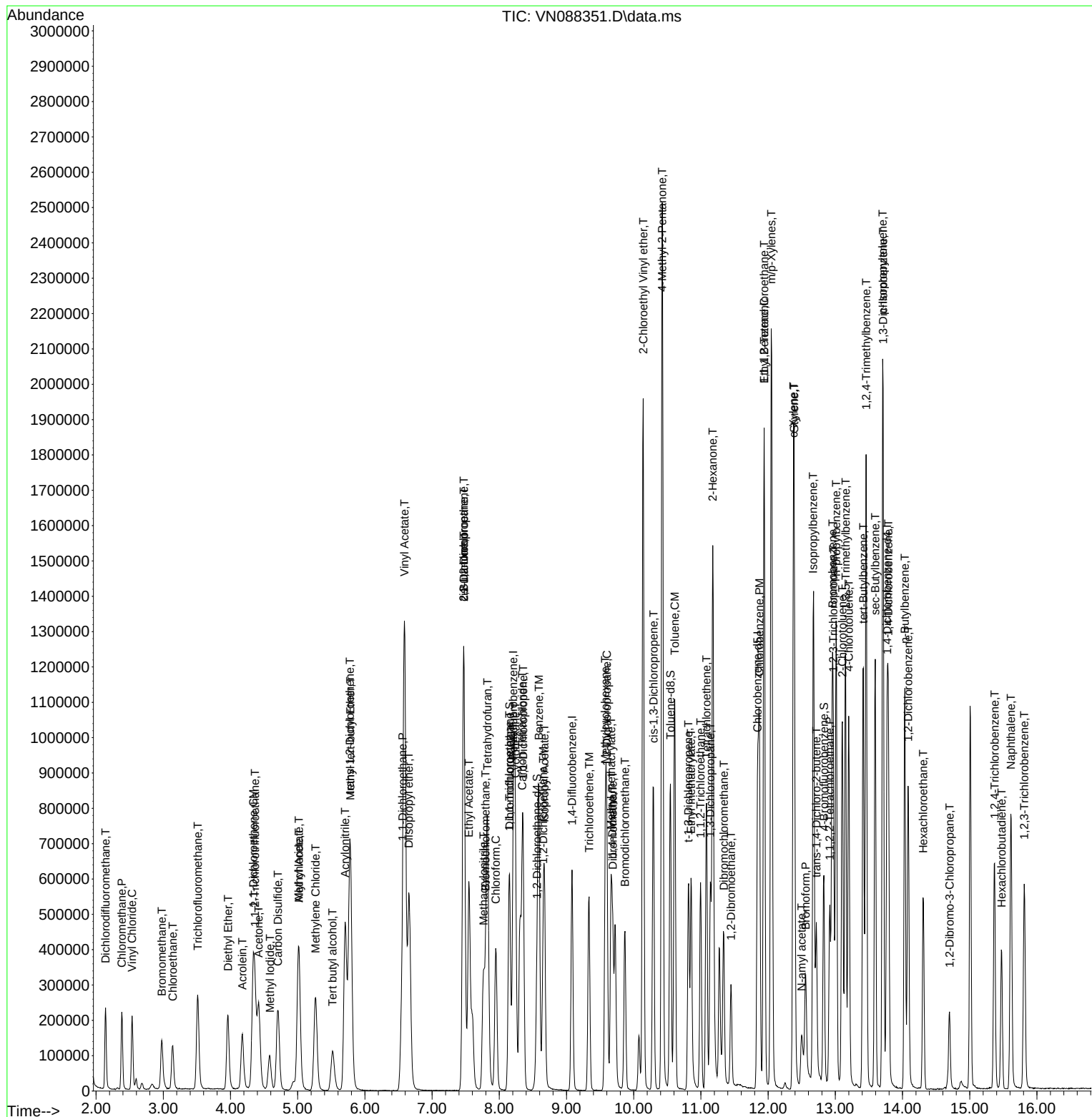
Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
----------	------	------	----------	------	-------	----------

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

Instrument :
MSVOA_N
ClientSampleId :
ICVVN112425

Manual Integrations APPROVED

Reviewed By :John Carlone 11/25/2025
Supervised By :Mahesh Dadoda 11/26/2025



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN112425\
 Data File : VN088351.D
 Acq On : 24 Nov 2025 15:18
 Operator : JC\MD
 Sample : VSTDICV050
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 ICSVN112425

Quant Time: Nov 25 01:39:30 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N112425W.M
 Quant Title : SW846 8260
 QLast Update : Tue Nov 25 01:34:52 2025
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
 Max. RRF Dev : 25% Max. Rel. Area : 150%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
1 I	Pentafluorobenzene	1.000	1.000	0.0	104	0.00
2 T	Dichlorodifluoromethane	0.751	0.772	-2.8	108	0.00
3 P	Chloromethane	0.807	0.848	-5.1	110	0.00
4 C	Vinyl Chloride	0.821	0.818	0.4#	105	0.00
5 T	Bromomethane	0.395	0.430	-8.9	113	0.00
6 T	Chloroethane	0.508	0.523	-3.0	108	0.00
7 T	Trichlorofluoromethane	1.153	1.149	0.3	110	0.00
8 T	Diethyl Ether	0.433	0.462	-6.7	108	0.00
9 T	1,1,2-Trichlorotrifluoroeth	0.614	0.578	5.9	107	0.00
10 T	Methyl Iodide	0.608	0.656	-7.9	103	0.00
11 T	Tert butyl alcohol	0.150	0.161	-7.3	104	0.00
12 CM	1,1-Dichloroethene	0.610	0.597	2.1#	109	0.00
13 T	Acrolein	0.144	0.154	-6.9	106	0.00
14 T	Allyl chloride	1.131	1.216	-7.5	110	0.00
15 T	Acrylonitrile	0.434	0.482	-11.1	107	0.00
16 T	Acetone	0.341	0.340	0.3	106	0.00
17 T	Carbon Disulfide	1.993	2.055	-3.1	108	0.00
18 T	Methyl Acetate	1.014	1.099	-8.4	109	0.00
19 T	Methyl tert-butyl Ether	2.285	2.624	-14.8	109	0.00
20 T	Methylene Chloride	0.729	0.789	-8.2	109	0.00
21 T	trans-1,2-Dichloroethene	0.676	0.722	-6.8	110	0.00
22 T	Diisopropyl ether	2.414	2.720	-12.7	110	0.00
23 T	Vinyl Acetate	1.945	2.203	-13.3	107	0.00
24 P	1,1-Dichloroethane	1.367	1.490	-9.0	109	0.00
25 T	2-Butanone	0.548	0.610	-11.3	107	0.00
26 T	2,2-Dichloropropane	1.180	1.273	-7.9	110	0.00
27 T	cis-1,2-Dichloroethene	0.785	0.846	-7.8	106	0.00
28 T	Bromochloromethane	0.681	0.634	6.9	90	0.00
29 T	Tetrahydrofuran	0.393	0.439	-11.7	106	0.00
30 C	Chloroform	1.343	1.450	-8.0#	108	0.00
31 T	Cyclohexane	1.239	1.194	3.6	110	0.00
32 T	1,1,1-Trichloroethane	1.181	1.208	-2.3	109	0.00
33 S	1,2-Dichloroethane-d4	0.942	0.887	5.8	92	0.00
34 I	1,4-Difluorobenzene	1.000	1.000	0.0	105	0.00
35 S	Dibromofluoromethane	0.346	0.303	12.4	91	0.00
36 T	1,1-Dichloropropene	0.535	0.502	6.2	107	0.00
37 T	Ethyl Acetate	0.703	0.723	-2.8	105	0.00
38 T	Carbon Tetrachloride	0.532	0.520	2.3	109	0.00
39 T	Methylcyclohexane	0.572	0.548	4.2	109	0.00
40 TM	Benzene	1.570	1.595	-1.6	107	0.00
41 T	Methacrylonitrile	0.348	0.375	-7.8	110	0.00
42 TM	1,2-Dichloroethane	0.599	0.632	-5.5	108	0.00
43 T	Isopropyl Acetate	1.074	1.142	-6.3	107	0.00
44 TM	Trichloroethene	0.371	0.367	1.1	108	0.00
45 C	1,2-Dichloropropane	0.402	0.419	-4.2#	108	0.00
46 T	Dibromomethane	0.287	0.302	-5.2	108	0.00
47 T	Bromodichloromethane	0.586	0.611	-4.3	109	0.00
48 T	Methyl methacrylate	0.490	0.535	-9.2	109	0.00

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN112425\
 Data File : VN088351.D
 Acq On : 24 Nov 2025 15:18
 Operator : JC\MD
 Sample : VSTDICV050
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 ICVVN112425

Quant Time: Nov 25 01:39:30 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N112425W.M
 Quant Title : SW846 8260
 QLast Update : Tue Nov 25 01:34:52 2025
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
 Max. RRF Dev : 25% Max. Rel. Area : 150%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
49 T	1,4-Dioxane	0.006	0.006	0.0	101	0.00
50 S	Toluene-d8	1.289	1.140	11.6	93	0.00
51 T	4-Methyl-2-Pentanone	0.633	0.693	-9.5	108	0.00
52 CM	Toluene	0.906	0.942	-4.0#	108	0.00
53 T	t-1,3-Dichloropropene	0.601	0.642	-6.8	107	0.00
54 T	cis-1,3-Dichloropropene	0.632	0.684	-8.2	108	0.00
55 T	1,1,2-Trichloroethane	0.351	0.361	-2.8	107	0.00
56 T	Ethyl methacrylate	0.554	0.642	-15.9	109	0.00
57 T	1,3-Dichloropropane	0.633	0.672	-6.2	109	0.00
58 T	2-Chloroethyl Vinyl ether	0.319	0.338	-6.0	97	0.00
59 T	2-Hexanone	0.399	0.444	-11.3	107	0.00
60 T	Dibromochloromethane	0.400	0.421	-5.2	108	0.00
61 T	1,2-Dibromoethane	0.355	0.380	-7.0	108	0.00
62 S	4-Bromofluorobenzene	0.442	0.397	10.2	93	0.00
63 I	Chlorobenzene-d5	1.000	1.000	0.0	104	0.00
64 T	Tetrachloroethene	0.396	0.370	6.6	102	0.00
65 PM	Chlorobenzene	1.123	1.156	-2.9	109	0.00
66 T	1,1,1,2-Tetrachloroethane	0.379	0.393	-3.7	107	0.00
67 C	Ethyl Benzene	1.963	2.027	-3.3#	109	0.00
68 T	m/p-Xylenes	0.728	0.766	-5.2	109	0.00
69 T	o-Xylene	0.694	0.727	-4.8	109	0.00
70 T	Styrene	1.127	1.240	-10.0	108	0.00
71 P	Bromoform	0.277	0.299	-7.9	109	0.00
72 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	102	0.00
73 T	Isopropylbenzene	3.697	3.909	-5.7	110	0.00
74 T	N-ethyl acetate	1.259	1.044	17.1	101	0.00
75 P	1,1,2,2-Tetrachloroethane	1.195	1.199	-0.3	108	0.00
76 T	1,2,3-Trichloropropane	1.171	1.271	-8.5	107	0.00
77 T	Bromobenzene	0.846	0.903	-6.7	106	0.00
78 T	n-propylbenzene	4.523	4.744	-4.9	109	0.00
79 T	2-Chlorotoluene	2.762	2.828	-2.4	106	0.00
80 T	1,3,5-Trimethylbenzene	3.063	3.243	-5.9	108	0.00
81 T	trans-1,4-Dichloro-2-butene	0.413	0.455	-10.2	107	0.00
82 T	4-Chlorotoluene	2.743	2.968	-8.2	106	0.00
83 T	tert-Butylbenzene	2.604	2.691	-3.3	109	0.00
84 T	1,2,4-Trimethylbenzene	3.045	3.320	-9.0	108	0.00
85 T	sec-Butylbenzene	3.741	3.837	-2.6	108	0.00
86 T	p-Isopropyltoluene	3.068	3.204	-4.4	107	0.00
87 T	1,3-Dichlorobenzene	1.647	1.689	-2.6	107	0.00
88 T	1,4-Dichlorobenzene	1.725	1.758	-1.9	106	0.00
89 T	n-Butylbenzene	2.979	3.041	-2.1	107	0.00
90 T	Hexachloroethane	0.560	0.558	0.4	106	0.00
91 T	1,2-Dichlorobenzene	1.550	1.625	-4.8	108	0.00
92 T	1,2-Dibromo-3-Chloropropane	0.290	0.294	-1.4	103	0.00
93 T	1,2,4-Trichlorobenzene	0.921	0.945	-2.6	108	0.00
94 T	Hexachlorobutadiene	0.331	0.304	8.2	104	0.00
95 T	Naphthalene	3.190	3.398	-6.5	105	0.00

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN112425\
Data File : VN088351.D
Acq On : 24 Nov 2025 15:18
Operator : JC\MD
Sample : VSTDICV050
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
ICVVN112425

Quant Time: Nov 25 01:39:30 2025
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N112425W.M
Quant Title : SW846 8260
QLast Update : Tue Nov 25 01:34:52 2025
Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
Max. RRF Dev : 25% Max. Rel. Area : 150%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
96 T	1,2,3-Trichlorobenzene	0.896	0.889	0.8	103	0.00

(#) = Out of Range

SPCC's out = 0 CCC's out = 6

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN112425\
 Data File : VN088351.D
 Acq On : 24 Nov 2025 15:18
 Operator : JC\MD
 Sample : VSTDICV050
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 ICSVN112425

Quant Time: Nov 25 01:39:30 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N112425W.M
 Quant Title : SW846 8260
 QLast Update : Tue Nov 25 01:34:52 2025
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
 Max. RRF Dev : 25% Max. Rel. Area : 150%

	Compound	Amount	Calc.	%Dev	Area%	Dev(min)
1 I	Pentafluorobenzene	50.000	50.000	0.0	104	0.00
2 T	Dichlorodifluoromethane	50.000	51.389	-2.8	108	0.00
3 P	Chloromethane	50.000	52.578	-5.2	110	0.00
4 C	Vinyl Chloride	50.000	49.842	0.3#	105	0.00
5 T	Bromomethane	50.000	54.393	-8.8	113	0.00
6 T	Chloroethane	50.000	51.450	-2.9	108	0.00
7 T	Trichlorofluoromethane	50.000	49.826	0.3	110	0.00
8 T	Diethyl Ether	50.000	53.384	-6.8	108	0.00
9 T	1,1,2-Trichlorotrifluoroeth	50.000	47.094	5.8	107	0.00
10 T	Methyl Iodide	50.000	53.976	-8.0	103	0.00
11 T	Tert butyl alcohol	250.000	268.272	-7.3	104	0.00
12 CM	1,1-Dichloroethene	50.000	48.957	2.1#	109	0.00
13 T	Acrolein	250.000	266.656	-6.7	106	0.00
14 T	Allyl chloride	50.000	53.793	-7.6	110	0.00
15 T	Acrylonitrile	250.000	277.614	-11.0	107	0.00
16 T	Acetone	250.000	249.797	0.1	106	0.00
17 T	Carbon Disulfide	50.000	51.558	-3.1	108	0.00
18 T	Methyl Acetate	50.000	54.196	-8.4	109	0.00
19 T	Methyl tert-butyl Ether	50.000	57.401	-14.8	109	0.00
20 T	Methylene Chloride	50.000	54.158	-8.3	109	0.00
21 T	trans-1,2-Dichloroethene	50.000	53.421	-6.8	110	0.00
22 T	Diisopropyl ether	50.000	56.335	-12.7	110	0.00
23 T	Vinyl Acetate	250.000	283.141	-13.3	107	0.00
24 P	1,1-Dichloroethane	50.000	54.516	-9.0	109	0.00
25 T	2-Butanone	250.000	278.613	-11.4	107	0.00
26 T	2,2-Dichloropropane	50.000	53.957	-7.9	110	0.00
27 T	cis-1,2-Dichloroethene	50.000	53.866	-7.7	106	0.00
28 T	Bromochloromethane	50.000	46.526	6.9	90	0.00
29 T	Tetrahydrofuran	250.000	279.515	-11.8	106	0.00
30 C	Chloroform	50.000	53.957	-7.9#	108	0.00
31 T	Cyclohexane	50.000	48.211	3.6	110	0.00
32 T	1,1,1-Trichloroethane	50.000	51.150	-2.3	109	0.00
33 S	1,2-Dichloroethane-d4	50.000	47.098	5.8	92	0.00
34 I	1,4-Difluorobenzene	50.000	50.000	0.0	105	0.00
35 S	Dibromofluoromethane	50.000	43.757	12.5	91	0.00
36 T	1,1-Dichloropropene	50.000	46.967	6.1	107	0.00
37 T	Ethyl Acetate	50.000	51.431	-2.9	105	0.00
38 T	Carbon Tetrachloride	50.000	48.837	2.3	109	0.00
39 T	Methylcyclohexane	50.000	47.913	4.2	109	0.00
40 TM	Benzene	50.000	50.800	-1.6	107	0.00
41 T	Methacrylonitrile	50.000	53.883	-7.8	110	0.00
42 TM	1,2-Dichloroethane	50.000	52.806	-5.6	108	0.00
43 T	Isopropyl Acetate	50.000	53.165	-6.3	107	0.00
44 TM	Trichloroethene	50.000	49.444	1.1	108	0.00
45 C	1,2-Dichloropropane	50.000	52.095	-4.2#	108	0.00
46 T	Dibromomethane	50.000	52.686	-5.4	108	0.00
47 T	Bromodichloromethane	50.000	52.155	-4.3	109	0.00
48 T	Methyl methacrylate	50.000	54.535	-9.1	109	0.00

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN112425\
 Data File : VN088351.D
 Acq On : 24 Nov 2025 15:18
 Operator : JC\MD
 Sample : VSTDICV050
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 ICVVN112425

Quant Time: Nov 25 01:39:30 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N112425W.M
 Quant Title : SW846 8260
 QLast Update : Tue Nov 25 01:34:52 2025
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
 Max. RRF Dev : 25% Max. Rel. Area : 150%

	Compound	Amount	Calc.	%Dev	Area%	Dev(min)
49 T	1,4-Dioxane	1000.000	1021.452	-2.1	101	0.00
50 S	Toluene-d8	50.000	44.193	11.6	93	0.00
51 T	4-Methyl-2-Pentanone	250.000	273.447	-9.4	108	0.00
52 CM	Toluene	50.000	51.998	-4.0#	108	0.00
53 T	t-1,3-Dichloropropene	50.000	53.428	-6.9	107	0.00
54 T	cis-1,3-Dichloropropene	50.000	54.081	-8.2	108	0.00
55 T	1,1,2-Trichloroethane	50.000	51.530	-3.1	107	0.00
56 T	Ethyl methacrylate	50.000	57.977	-16.0	109	0.00
57 T	1,3-Dichloropropane	50.000	53.102	-6.2	109	0.00
58 T	2-Chloroethyl Vinyl ether	250.000	264.613	-5.8	97	0.00
59 T	2-Hexanone	250.000	278.086	-11.2	107	0.00
60 T	Dibromochloromethane	50.000	52.650	-5.3	108	0.00
61 T	1,2-Dibromoethane	50.000	53.546	-7.1	108	0.00
62 S	4-Bromofluorobenzene	50.000	44.823	10.4	93	0.00
63 I	Chlorobenzene-d5	50.000	50.000	0.0	104	0.00
64 T	Tetrachloroethene	50.000	46.666	6.7	102	0.00
65 PM	Chlorobenzene	50.000	51.452	-2.9	109	0.00
66 T	1,1,1,2-Tetrachloroethane	50.000	51.838	-3.7	107	0.00
67 C	Ethyl Benzene	50.000	51.635	-3.3#	109	0.00
68 T	m/p-Xylenes	100.000	105.191	-5.2	109	0.00
69 T	o-Xylene	50.000	52.376	-4.8	109	0.00
70 T	Styrene	50.000	55.031	-10.1	108	0.00
71 P	Bromoform	50.000	53.811	-7.6	109	0.00
72 I	1,4-Dichlorobenzene-d4	50.000	50.000	0.0	102	0.00
73 T	Isopropylbenzene	50.000	52.866	-5.7	110	0.00
74 T	N-ethyl acetate	50.000	41.449	17.1	101	0.00
75 P	1,1,2,2-Tetrachloroethane	50.000	50.155	-0.3	108	0.00
76 T	1,2,3-Trichloropropane	50.000	54.274	-8.5	107	0.00
77 T	Bromobenzene	50.000	53.406	-6.8	106	0.00
78 T	n-propylbenzene	50.000	52.446	-4.9	109	0.00
79 T	2-Chlorotoluene	50.000	51.202	-2.4	106	0.00
80 T	1,3,5-Trimethylbenzene	50.000	52.936	-5.9	108	0.00
81 T	trans-1,4-Dichloro-2-butene	50.000	55.117	-10.2	107	0.00
82 T	4-Chlorotoluene	50.000	54.093	-8.2	106	0.00
83 T	tert-Butylbenzene	50.000	51.660	-3.3	109	0.00
84 T	1,2,4-Trimethylbenzene	50.000	54.519	-9.0	108	0.00
85 T	sec-Butylbenzene	50.000	51.285	-2.6	108	0.00
86 T	p-Isopropyltoluene	50.000	52.216	-4.4	107	0.00
87 T	1,3-Dichlorobenzene	50.000	51.282	-2.6	107	0.00
88 T	1,4-Dichlorobenzene	50.000	50.949	-1.9	106	0.00
89 T	n-Butylbenzene	50.000	51.040	-2.1	107	0.00
90 T	Hexachloroethane	50.000	49.768	0.5	106	0.00
91 T	1,2-Dichlorobenzene	50.000	52.435	-4.9	108	0.00
92 T	1,2-Dibromo-3-Chloropropane	50.000	50.621	-1.2	103	0.00
93 T	1,2,4-Trichlorobenzene	50.000	51.301	-2.6	108	0.00
94 T	Hexachlorobutadiene	50.000	45.969	8.1	104	0.00
95 T	Naphthalene	50.000	53.258	-6.5	105	0.00

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN112425\
Data File : VN088351.D
Acq On : 24 Nov 2025 15:18
Operator : JC\MD
Sample : VSTDICV050
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
ICVVN112425

Quant Time: Nov 25 01:39:30 2025
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N112425W.M
Quant Title : SW846 8260
QLast Update : Tue Nov 25 01:34:52 2025
Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
Max. RRF Dev : 25% Max. Rel. Area : 150%

	Compound	Amount	Calc.	%Dev	Area%	Dev(min)
96 T	1,2,3-Trichlorobenzene	50.000	49.620	0.8	103	0.00

(#) = Out of Range SPCC's out = 0 CCC's out = 6



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

VOLATILE ORGANICS INITIAL CALIBRATION DATA

Lab Name: Alliance Contract: JAC005
 Lab Code: ACE SDG No.: Q3757
 Instrument ID: MSVOA_X Calibration Date(s): 12/04/2025 12/04/2025
 Heated Purge: (Y/N) N Calibration Time(s): 08:47 10:38
 GC Column: DB-624UI ID: 0.18 (mm)

LAB FILE ID:		RRF001 = VX048714.D		RRF005 = VX048715.D		RRF020 = VX048716.D			
		RRF050 = VX048717.D		RRF100 = VX048718.D		RRF150 = VX048719.D			
COMPOUND	RRF001	RRF005	RRF020	RRF050	RRF100	RRF150	RRF	% RSD	
Vinyl Chloride	0.704	0.624	0.667	0.709	0.700	0.704	0.685	4.9	
1,1-Dichloroethene	0.607	0.533	0.560	0.601	0.584	0.598	0.581	4.9	
1,1-Dichloroethane	1.208	0.803	1.038	0.910	1.223	0.914	1.016	16.9	
cis-1,2-Dichloroethene	0.776	0.559	0.634	0.629	0.785	0.630	0.669	13.6	
1,1,1-Trichloroethane	1.091	0.859	0.955	0.912	1.013	1.023	0.975	8.6	
Benzene	1.508	1.396	1.314	1.323	1.402	1.444	1.398	5.2	
1,2-Dichloroethane	0.554	0.497	0.439	0.452	0.525	0.502	0.495	8.7	
Trichloroethene	0.387	0.340	0.363	0.367	0.345	0.387	0.365	5.4	
1,1,2-Trichloroethane	0.359	0.316	0.325	0.378	0.310	0.347	0.339	7.9	
Tetrachloroethene	0.383	0.367	0.358	0.353	0.315	0.410	0.364	8.8	
1,2-Dichloroethane-d4		0.677	0.618	0.578	0.787	0.695	0.671	11.9	
Dibromofluoromethane		0.316	0.359	0.334	0.355	0.357	0.344	5.4	
Toluene-d8		1.281	1.161	1.312	1.051	1.228	1.207	8.6	
4-Bromofluorobenzene		0.416	0.418	0.459	0.397	0.391	0.416	6.4	

* 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_X\Method\
 Method File : 82X120425W.M
 Title : SW846 8260
 Last Update : Fri Dec 05 03:27:34 2025
 Response Via : Initial Calibration

Calibration Files

1 =VX048714.D 5 =VX048715.D 20 =VX048716.D 50 =VX048717.D 100 =VX048718.D 150 =VX048719.D

	Compound	1	5	20	50	100	150	Avg	%RSD

1) I	Pentafluorobenzene	-----ISTD-----							
2) T	Dichlorodifluo...	0.478	0.435	0.536	0.572	0.496	0.540	0.509	9.72
3) P	Chloromethane	0.725	0.555	0.597	0.645	0.671	0.664	0.643	9.29
4) C	Vinyl Chloride	0.704	0.624	0.667	0.709	0.700	0.704	0.685	4.87#
5) T	Bromomethane		0.439	0.431	0.467	0.441	0.338	0.423	11.68
6) T	Chloroethane	0.516	0.366	0.394	0.426	0.434	0.420	0.426	11.88
7) T	Trichlorofluor...	0.999	0.934	0.972	1.032	0.940	1.011	0.981	4.01
8) T	Diethyl Ether	0.445	0.357	0.341	0.380	0.435	0.392	0.392	10.58
9) T	1,1,2-Trichlor...	0.591	0.565	0.555	0.596	0.522	0.585	0.569	4.90
10) T	Methyl Iodide		0.589	0.745	0.865	0.957	0.893	0.810	17.96
11) T	Tert butyl alc...		0.086	0.106	0.115	0.137	0.098	0.108	17.56
12) CM	1,1-Dichloroet...	0.607	0.533	0.560	0.601	0.584	0.598	0.581	4.95#
13) T	Acrolein		0.065	0.067	0.086	0.117	0.119	0.091	28.91
14) T	Allyl chloride	1.104	0.725	0.973	1.075	1.172	0.799	0.975	18.29
15) T	Acrylonitrile	0.366	0.256	0.326	0.367	0.423	0.288	0.338	17.97
16) T	Acetone	0.319	0.213	0.221	0.247	0.286	0.255	0.257	15.61
17) T	Carbon Disulfide	1.975	1.314	1.640	1.773	1.732	1.683	1.686	12.83
18) T	Methyl Acetate	0.831	0.525	0.697	0.777	0.907	0.600	0.723	19.86
19) T	Methyl tert-bu...	2.162	1.452	1.778	1.996	2.253	1.697	1.890	16.03
20) T	Methylene Chlo...	0.841	0.524	0.586	0.650	0.718	0.534	0.642	18.95
21) T	trans-1,2-Dich...	0.705	0.487	0.571	0.626	0.642	0.524	0.593	13.63
22) T	Diisopropyl ether	2.158	1.457	1.895	1.567	2.280	1.573	1.822	18.81
23) T	Vinyl Acetate	1.537	1.176	1.611	1.382	2.071	1.407	1.531	19.86
24) P	1,1-Dichloroet...	1.208	0.803	1.038	0.910	1.223	0.914	1.016	16.91
25) T	2-Butanone	0.399	0.286	0.402	0.337	0.507	0.344	0.379	20.00
26) T	2,2-Dichloropr...	1.068	0.780	0.883	0.831	0.994	0.839	0.899	12.22
27) T	cis-1,2-Dichlo...	0.776	0.559	0.634	0.629	0.785	0.630	0.669	13.63
28) T	Bromochloromet...	0.565	0.416	0.496	0.417	0.611	0.483	0.498	15.73
29) T	Tetrahydrofuran	0.318	0.214	0.287	0.248	0.363	0.306	0.290	18.22
30) C	Chloroform	1.260	0.905	1.053	0.981	1.212	1.121	1.089	12.46#
31) T	Cyclohexane		0.969	0.963	0.797	0.861	0.946	0.907	8.33
32) T	1,1,1-Trichlor...	1.091	0.859	0.955	0.912	1.013	1.023	0.975	8.59
33) S	1,2-Dichloroet...		0.677	0.618	0.578	0.787	0.695	0.671	11.91
34) I	1,4-Difluorobenzene	-----ISTD-----							
35) S	Dibromofluorom...		0.316	0.359	0.334	0.355	0.357	0.344	5.37
36) T	1,1-Dichloropr...	0.522	0.461	0.516	0.439	0.422	0.469	0.472	8.63
37) T	Ethyl Acetate	0.555	0.474	0.655	0.540	0.688	0.469	0.563	16.19
38) T	Carbon Tetrach...	0.501	0.572	0.590	0.528	0.476	0.538	0.534	8.00
39) T	Methylcyclohexane	0.539	0.540	0.550	0.642	0.478	0.596	0.558	10.06
40) TM	Benzene	1.508	1.396	1.314	1.323	1.402	1.444	1.398	5.24
41) T	Methacrylonitrile	0.217	0.205	0.302	0.242	0.307	0.264	0.256	16.65
42) TM	1,2-Dichloroet...	0.554	0.497	0.439	0.452	0.525	0.502	0.495	8.74
43) T	Isopropyl Acetate	0.878	0.867	0.738	0.756	0.969	0.910	0.853	10.49
44) TM	Trichloroethene	0.387	0.340	0.363	0.367	0.345	0.387	0.365	5.42
45) C	1,2-Dichloropr...	0.339	0.341	0.314	0.383	0.360	0.360	0.349	6.75#
46) T	Dibromomethane	0.256	0.257	0.242	0.289	0.273	0.271	0.265	6.18
47) T	Bromodichlorom...	0.595	0.538	0.488	0.587	0.555	0.551	0.552	6.98
48) T	Methyl methacr...	0.427	0.395	0.355	0.465	0.480	0.453	0.429	10.99
49) T	1,4-Dioxane	0.006	0.005	0.005	0.006	0.007	0.006	0.006	10.25
50) S	Toluene-d8		1.281	1.161	1.312	1.051	1.228	1.207	8.63
51) T	4-Methyl-2-Pen...	0.547	0.544	0.470	0.605	0.460	0.552	0.530	10.40
52) CM	Toluene	0.838	0.786	0.831	0.964	0.751	0.888	0.843	8.96#
53) T	t-1,3-Dichloro...	0.543	0.445	0.490	0.600	0.495	0.581	0.526	11.30
54) T	cis-1,3-Dichlo...	0.540	0.537	0.531	0.644	0.524	0.613	0.565	8.94
55) T	1,1,2-Trichlor...	0.359	0.316	0.325	0.378	0.310	0.347	0.339	7.85
56) T	Ethyl methacry...	0.492	0.440	0.502	0.627	0.520	0.603	0.531	13.33

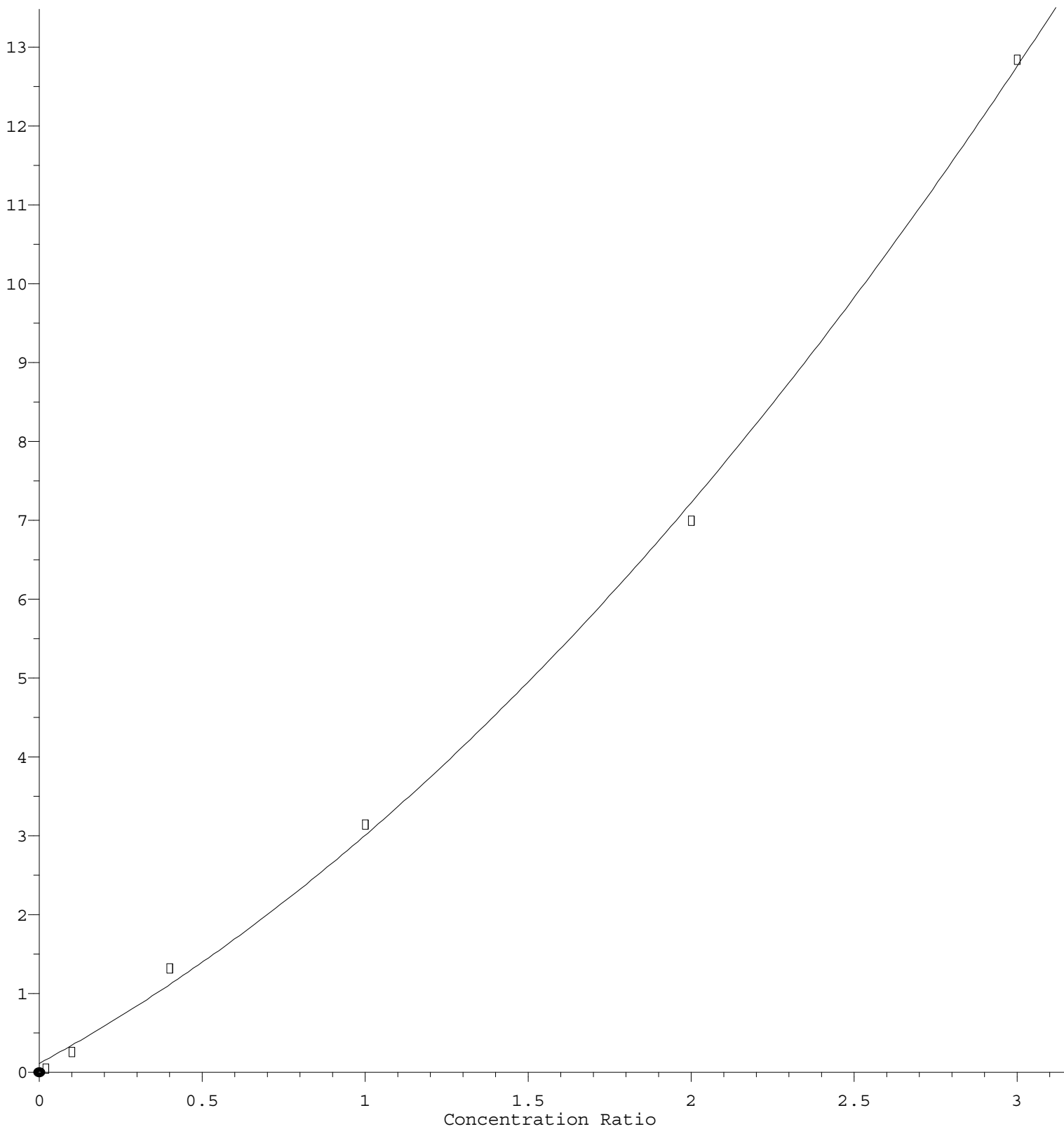
Method Path : Z:\voasrv\HPCHEM1\MSVOA_X\Method\
 Method File : 82X120425W.M

57)	T	1,3-Dichloropr...	0.620	0.509	0.535	0.628	0.513	0.579	0.564	9.39
58)	T	2-Chloroethyl ...	0.227	0.295	0.268	0.344	0.256	0.332	0.287	15.76
59)	T	2-Hexanone	0.383	0.311	0.345	0.442	0.342	0.408	0.372	13.02
60)	T	Dibromochlorom...	0.415	0.390	0.412	0.467	0.401	0.431	0.419	6.54
61)	T	1,2-Dibromoethane	0.388	0.334	0.356	0.411	0.343	0.372	0.367	7.85
62)	S	4-Bromofluorob...	0.416	0.418	0.459	0.397	0.391	0.416		6.39
63)	I	Chlorobenzene-d5	-----ISTD-----							
64)	T	Tetrachloroethene	0.383	0.367	0.358	0.353	0.315	0.410	0.364	8.76
65)	PM	Chlorobenzene	1.181	1.082	1.060	1.087	1.044	1.113	1.094	4.45
66)	T	1,1,1,2-Tetrac...	0.434	0.376	0.367	0.380	0.382	0.410	0.391	6.47
67)	C	Ethyl Benzene	1.741	1.720	1.759	1.893	1.702	1.900	1.786	4.93#
68)	T	m/p-Xylenes	0.699	0.675	0.684	0.717	0.664	0.736	0.696	3.89
69)	T	o-Xylene	0.577	0.634	0.658	0.678	0.653	0.712	0.652	6.98
70)	T	Styrene	0.989	1.068	1.129	1.170	1.134	1.223	1.119	7.28
71)	P	Bromoform	0.334	0.331	0.335	0.328	0.372	0.395	0.349	7.94
72)	I	1,4-Dichlorobenzen...	-----ISTD-----							
73)	T	Isopropylbenzene	3.175	3.265	3.404	3.633	3.155	3.586	3.369	6.12
74)	T	N-amyl acetate	1.397	1.236	1.349	1.738	1.491	1.572	1.464	12.11
75)	P	1,1,2,2-Tetrac...	1.297	1.035	1.035	1.179	1.108	1.168	1.137	8.78
76)	T	1,2,3-Trichlor...	1.093	0.906	0.839	0.958	0.885	0.906	0.931	9.43
77)	T	Bromobenzene	0.906	0.870	0.864	0.886	0.905	0.965	0.899	4.06
78)	T	n-propylbenzene	3.763	3.672	3.959	4.306	3.582	4.133	3.902	7.19
79)	T	2-Chlorotoluene	2.399	2.263	2.342	2.521	2.260	2.493	2.380	4.70
80)	T	1,3,5-Trimethy...	2.581	2.644	2.831	2.948	2.635	2.978	2.770	6.22
81)	T	trans-1,4-Dich...	0.318	0.364	0.416	0.416	0.447	0.392		12.98
82)	T	4-Chlorotoluene	2.746	2.550	2.779	2.984	2.667	2.936	2.777	5.87
83)	T	tert-Butylbenzene	2.630	2.796	2.928	3.026	2.744	3.143	2.878	6.61
84)	T	1,2,4-Trimethy...	2.583	2.631	2.811	2.983	2.715	3.039	2.794	6.66
85)	T	sec-Butylbenzene	3.257	3.441	3.572	3.766	3.161	3.728	3.488	7.08
86)	T	p-Isopropyltol...	2.719	2.915	3.051	3.180	2.786	3.236	2.981	7.05
87)	T	1,3-Dichlorobe...	1.691	1.611	1.630	1.647	1.612	1.756	1.658	3.41
88)	T	1,4-Dichlorobe...	1.952	1.640	1.612	1.662	1.639	1.775	1.713	7.60
89)	T	n-Butylbenzene	2.735	2.610	3.077	2.959	2.457	2.909	2.791	8.34
90)	T	Hexachloroethane	0.618	0.555	0.627	0.594	0.538	0.718	0.608	10.50
91)	T	1,2-Dichlorobe...	1.632	1.491	1.702	1.582	1.557	1.680	1.607	4.95
92)	T	1,2-Dibromo-3-...	0.293	0.230	0.273	0.268	0.264	0.332	0.277	12.28
93)	T	1,2,4-Trichlor...	1.053	0.919	1.043	0.997	1.062	1.295	1.062	11.87
94)	T	Hexachlorobuta...	0.500	0.437	0.452	0.395	0.390	0.460	0.439	9.52
95)	T	Naphthalene	2.336	2.578	3.292	3.142	3.497	4.280	3.187	21.73
96)	T	1,2,3-Trichlor...	0.912	0.836	0.989	0.930	1.005	1.203	0.979	12.77

(#) = Out of Range

Naphthalene

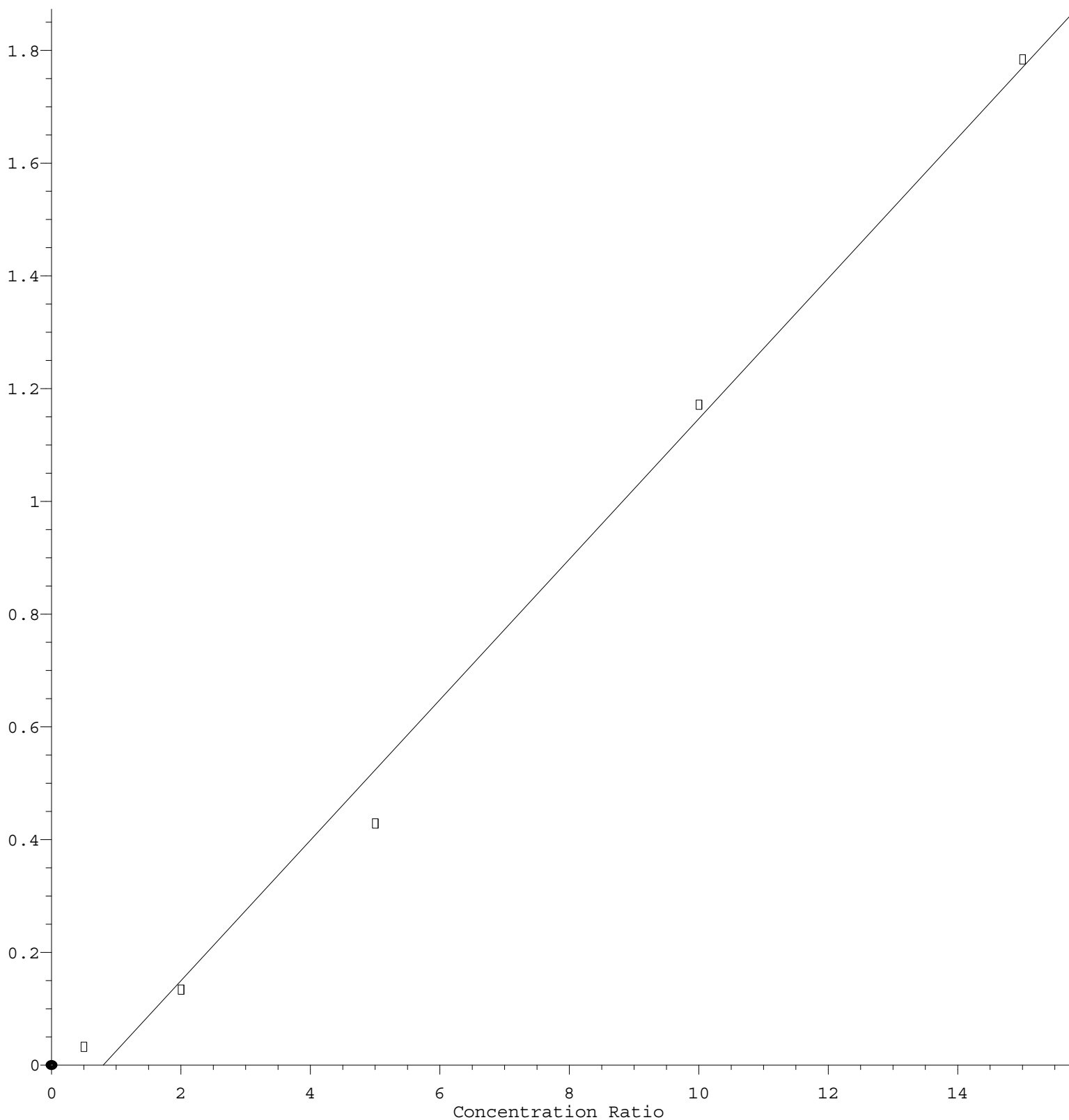
Response Ratio



$R = 6.609e-001 A^2 + 2.232e+000 A + 1.131e-001$
 Coef of Det (r^2) = 0.998889 Curve Fit: Quadratic
 Method Name: Z:\voasrv\HPCHEM1\MSVOA X\Method\82X120425W.M
 Calibration Table Last Updated: Fri Dec 05 03:27:34 2025

Acrolein

Response Ratio



$\text{Response} = 1.246\text{e-}001 * \text{Amt} - 1.002\text{e-}001$
 Coef of Det (r^2) = 0.993311 Curve Fit: Linear
 Method Name: Z:\voasrv\HPCHEM1\MSVOA X\Method\82X120425W.M
 Calibration Table Last Updated: Fri Dec 05 03:27:34 2025

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120425\
 Data File : VX048714.D
 Acq On : 04 Dec 2025 08:47
 Operator : JC/MD
 Sample : VSTDICC001
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 VSTDICC001

Manual Integrations
 APPROVED

Reviewed By : John Carlone 12/05/2025
 Supervised By : Semsettin Yesilyurt 12/05/2025

Quant Time: Dec 05 03:18:22 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X120425W.M
 Quant Title : SW846 8260
 QLast Update : Fri Dec 05 03:18:21 2025
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.531	168	206937	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.739	114	356626	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.037	117	323968	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.006	152	158289	50.000	ug/l	0.00

System Monitoring Compounds

33) 1,2-Dichloroethane-d4	0.000	65	0d	0.000	ug/l	
Spiked Amount	50.000	Range	78 - 117	Recovery	=	0.000%#
35) Dibromofluoromethane	0.000	113	0d	0.000	ug/l	
Spiked Amount	50.000	Range	75 - 124	Recovery	=	0.000%#
50) Toluene-d8	0.000	98	0d	0.000	ug/l	
Spiked Amount	50.000	Range	92 - 112	Recovery	=	0.000%#
62) 4-Bromofluorobenzene	0.000	95	0d	0.000	ug/l	
Spiked Amount	50.000	Range	83 - 123	Recovery	=	0.000%#

Target Compounds

					Qvalue	
2) Dichlorodifluoromethane	1.172	85	1978	0.938	ug/l #	66
3) Chloromethane	1.300	50	3002	1.128	ug/l	92
4) Vinyl Chloride	1.380	62	2913	1.028	ug/l	99
6) Chloroethane	1.697	64	2136	1.211	ug/l	98
7) Trichlorofluoromethane	1.898	101	4135	1.018	ug/l	93
8) Diethyl Ether	2.130	74	1841	1.136	ug/l	91
9) 1,1,2-Trichlorotrifluo...	2.343	101	2444	1.038	ug/l	95
12) 1,1-Dichloroethene	2.325	96	2514	1.046	ug/l	93
14) Allyl chloride	2.660	41	4571	1.133	ug/l	94
15) Acrylonitrile	3.056	53	7576	5.422	ug/l	95
16) Acetone	2.373	43	6606	6.213	ug/l	94
17) Carbon Disulfide	2.514	76	8174	1.171	ug/l	96
18) Methyl Acetate	2.703	43	3438	1.149	ug/l	95
19) Methyl tert-butyl Ether	3.105	73	8947	1.144	ug/l #	90
20) Methylene Chloride	2.788	84	3480	1.309	ug/l	88
21) trans-1,2-Dichloroethene	3.093	96	2919	1.190	ug/l	82
22) Diisopropyl ether	3.745	45	8931	1.184	ug/l #	88
23) Vinyl Acetate	3.721	43	31813	5.021	ug/l	98
24) 1,1-Dichloroethane	3.599	63	5001	1.189	ug/l	97
25) 2-Butanone	4.544	43	8255	5.260	ug/l	97
26) 2,2-Dichloropropane	4.458	77	4420	1.188	ug/l	99
27) cis-1,2-Dichloroethene	4.477	96	3211	1.160	ug/l	88
28) Bromochloromethane	4.879	49	2337	1.134	ug/l #	83
29) Tetrahydrofuran	4.989	42	6590	5.500	ug/l #	89
30) Chloroform	5.074	83	5213	1.157	ug/l	99
32) 1,1,1-Trichloroethane	5.361	97	4515	1.118	ug/l #	48
36) 1,1-Dichloropropene	5.678	75	3726	1.108	ug/l	91
37) Ethyl Acetate	4.714	43	3958m	0.985	ug/l	
38) Carbon Tetrachloride	5.659	117	3570	0.937	ug/l #	83
39) Methylcyclohexane	7.354	83	3845	0.967	ug/l #	87
40) Benzene	6.019	78	10753	1.079	ug/l	96
41) Methacrylonitrile	4.903	41	1547	0.846	ug/l #	62
42) 1,2-Dichloroethane	6.074	62	3949	1.119	ug/l	87
43) Isopropyl Acetate	6.318	43	6262	1.029	ug/l #	94
44) Trichloroethene	7.104	130	2758	1.060	ug/l	99
45) 1,2-Dichloropropane	7.409	63	2415	0.969	ug/l #	77

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120425\
 Data File : VX048714.D
 Acq On : 04 Dec 2025 08:47
 Operator : JC/MD
 Sample : VSTDICC001
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 VSTDICC001

Manual Integrations
 APPROVED

Reviewed By : John Carlone 12/05/2025
 Supervised By : Semsettin Yesilyurt 12/05/2025

Quant Time: Dec 05 03:18:22 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X120425W.M
 Quant Title : SW846 8260
 QLast Update : Fri Dec 05 03:18:21 2025
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
46) Dibromomethane	7.568	93	1828	0.967	ug/l	95
47) Bromodichloromethane	7.799	83	4244	1.077	ug/l #	85
48) Methyl methacrylate	7.677	41	3046	0.995	ug/l	93
49) 1,4-Dioxane	7.671	88	794	18.860	ug/l #	86
51) 4-Methyl-2-Pentanone	8.555	43	19490	5.160	ug/l	97
52) Toluene	8.702	92	5975	0.994	ug/l	95
53) t-1,3-Dichloropropene	8.970	75	3870	1.032	ug/l	99
54) cis-1,3-Dichloropropene	8.354	75	3853	0.956	ug/l #	93
55) 1,1,2-Trichloroethane	9.134	97	2561	1.059	ug/l #	90
56) Ethyl methacrylate	9.098	69	3510	0.927	ug/l	91
57) 1,3-Dichloropropane	9.293	76	4423	1.100	ug/l	97
58) 2-Chloroethyl Vinyl ether	8.232	63	8078	3.947	ug/l	97
59) 2-Hexanone	9.415	43	13642	5.145	ug/l	92
60) Dibromochloromethane	9.506	129	2961	0.990	ug/l	93
61) 1,2-Dibromoethane	9.592	107	2765	1.056	ug/l	97
64) Tetrachloroethene	9.256	164	2484	1.052	ug/l #	80
65) Chlorobenzene	10.061	112	7654	1.079	ug/l	99
66) 1,1,1,2-Tetrachloroethane	10.140	131	2813	1.109	ug/l #	61
67) Ethyl Benzene	10.177	91	11281	0.975	ug/l	95
68) m/p-Xylenes	10.281	106	9059	2.010	ug/l	99
69) o-Xylene	10.622	106	3736	0.885	ug/l	84
70) Styrene	10.640	104	6411	0.884	ug/l	96
71) Bromoform	10.787	173	2166	0.957	ug/l #	97
73) Isopropylbenzene	10.945	105	10050	0.942	ug/l	100
74) N-amyl acetate	10.829	43	4424	0.955	ug/l	95
75) 1,1,2,2-Tetrachloroethane	11.195	83	4105	1.141	ug/l	98
76) 1,2,3-Trichloropropane	11.219	75	3459m	1.139	ug/l	
77) Bromobenzene	11.183	156	2868	1.007	ug/l	95
78) n-propylbenzene	11.286	91	11914	0.964	ug/l	99
79) 2-Chlorotoluene	11.347	91	7596	1.008	ug/l	100
80) 1,3,5-Trimethylbenzene	11.433	105	8172	0.932	ug/l	96
82) 4-Chlorotoluene	11.439	91	8694	0.989	ug/l	99
83) tert-Butylbenzene	11.695	119	8327	0.914	ug/l	98
84) 1,2,4-Trimethylbenzene	11.732	105	8178	0.925	ug/l	99
85) sec-Butylbenzene	11.872	105	10311	0.934	ug/l	98
86) p-Isopropyltoluene	11.988	119	8607	0.912	ug/l	87
87) 1,3-Dichlorobenzene	11.951	146	5352	1.020	ug/l	99
88) 1,4-Dichlorobenzene	12.024	146	6180m	1.139	ug/l	
89) n-Butylbenzene	12.317	91	8657	0.980	ug/l	97
90) Hexachloroethane	12.524	117	1957	1.016	ug/l	90
91) 1,2-Dichlorobenzene	12.317	146	5165	1.015	ug/l	98
92) 1,2-Dibromo-3-Chloropr...	12.926	75	929	1.061	ug/l	74
93) 1,2,4-Trichlorobenzene	13.573	180	3332	0.991	ug/l	92
94) Hexachlorobutadiene	13.707	225	1584	1.139	ug/l	93
95) Naphthalene	13.762	128	7395	0.733	ug/l	99
96) 1,2,3-Trichlorobenzene	13.945	180	2886	0.931	ug/l	90

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

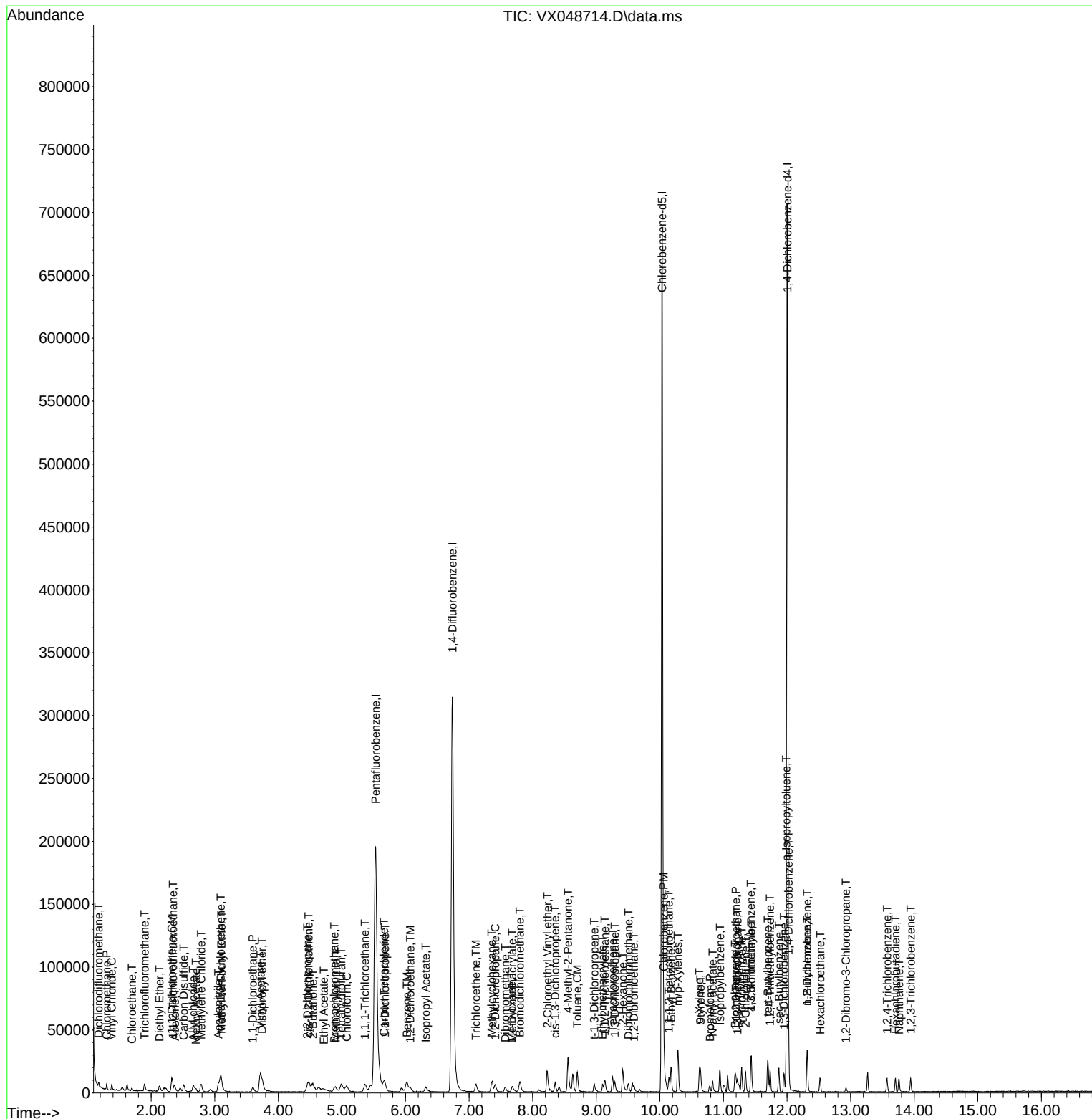
Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120425\
Data File : VX048714.D
Acq On : 04 Dec 2025 08:47
Operator : JC/MD
Sample : VSTDICC001
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 2 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
VSTDICC001

Manual Integrations
APPROVED

Reviewed By :John Carlone 12/05/2025
Supervised By :Semsettin Yesilyurt 12/05/2025

Quant Time: Dec 05 03:18:22 2025
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X120425W.M
Quant Title : SW846 8260
QLast Update : Fri Dec 05 03:18:21 2025
Response via : Initial Calibration



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120425\
 Data File : VX048715.D
 Acq On : 04 Dec 2025 09:16
 Operator : JC/MD
 Sample : VSTDICC005
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 VSTDICC005

Manual Integrations
 APPROVED

Reviewed By : John Carlone 12/05/2025
 Supervised By : Semsettin Yesilyurt 12/05/2025

Quant Time: Dec 05 02:07:22 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X120425W.M
 Quant Title : SW846 8260
 QLast Update : Fri Dec 05 02:05:18 2025
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.532	168	228640	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.739	114	358720	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.037	117	300438	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.006	152	155633	50.000	ug/l	0.00

System Monitoring Compounds

33) 1,2-Dichloroethane-d4	5.934	65	15478	5.044	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	10.080%#
35) Dibromofluoromethane	5.379	113	11348	4.594	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	9.180%#
50) Toluene-d8	8.629	98	45948	5.307	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	10.620%#
62) 4-Bromofluorobenzene	11.061	95	14926	4.999	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	10.000%#

Target Compounds

						Qvalue
2) Dichlorodifluoromethane	1.173	85	9943	4.268	ug/l	99
3) Chloromethane	1.301	50	12695	4.318	ug/l	98
4) Vinyl Chloride	1.380	62	14273	4.559	ug/l	93
5) Bromomethane	1.611	94	10033	5.186	ug/l	99
6) Chloroethane	1.691	64	8374	4.298	ug/l	98
7) Trichlorofluoromethane	1.892	101	21348	4.758	ug/l	99
8) Diethyl Ether	2.130	74	8173	4.563	ug/l	100
9) 1,1,2-Trichlorotrifluo...	2.337	101	12909	4.963	ug/l	97
10) Methyl Iodide	2.453	142	13462	3.635	ug/l	99
11) Tert butyl alcohol	2.941	59	9869	19.919	ug/l	99
12) 1,1-Dichloroethene	2.319	96	12191	4.592	ug/l	96
13) Acrolein	2.233	56	7426	53.206	ug/l	98
14) Allyl chloride	2.660	41	16575	3.718	ug/l	97
15) Acrylonitrile	3.056	53	29219	18.927	ug/l	91
16) Acetone	2.374	43	24360	20.736	ug/l	98
17) Carbon Disulfide	2.514	76	30036	3.896	ug/l	100
18) Methyl Acetate	2.697	43	12011m	3.634	ug/l	
19) Methyl tert-butyl Ether	3.099	73	33202	3.843	ug/l	95
20) Methylene Chloride	2.782	84	11979	4.079	ug/l	84
21) trans-1,2-Dichloroethene	3.087	96	11131	4.108	ug/l	96
22) Diisopropyl ether	3.745	45	33324	4.000	ug/l	97
23) Vinyl Acetate	3.709	43	134450	19.207	ug/l	99
24) 1,1-Dichloroethane	3.605	63	18354	3.951	ug/l	100
25) 2-Butanone	4.532	43	32730	18.877	ug/l	98
26) 2,2-Dichloropropane	4.459	77	17823	4.336	ug/l	98
27) cis-1,2-Dichloroethene	4.471	96	12772	4.177	ug/l	97
28) Bromochloromethane	4.867	49	9519	4.181	ug/l	91
29) Tetrahydrofuran	4.977	42	24477	18.488	ug/l	94
30) Chloroform	5.080	83	20698	4.158	ug/l	93
31) Cyclohexane	5.446	56	22164	5.344	ug/l	93
32) 1,1,1-Trichloroethane	5.361	97	19635	4.402	ug/l	93
36) 1,1-Dichloropropene	5.672	75	16539	4.889	ug/l	99
37) Ethyl Acetate	4.696	43	16990m	4.203	ug/l	
38) Carbon Tetrachloride	5.653	117	20529	5.357	ug/l	92
39) Methylcyclohexane	7.360	83	19359	4.840	ug/l	94
40) Benzene	6.019	78	50087	4.994	ug/l	98

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120425\
 Data File : VX048715.D
 Acq On : 04 Dec 2025 09:16
 Operator : JC/MD
 Sample : VSTDICC005
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 VSTDICC005

Manual Integrations
 APPROVED

Reviewed By :John Carlone 12/05/2025
 Supervised By :Semsettin Yesilyurt 12/05/2025

Quant Time: Dec 05 02:07:22 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X120425W.M
 Quant Title : SW846 8260
 QLast Update : Fri Dec 05 02:05:18 2025
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Methacrylonitrile	4.904	41	7366	4.005	ug/l	97
42) 1,2-Dichloroethane	6.068	62	17832	5.024	ug/l	97
43) Isopropyl Acetate	6.312	43	31107	5.082	ug/l	95
44) Trichloroethene	7.104	130	12207	4.665	ug/l	96
45) 1,2-Dichloropropane	7.409	63	12229	4.879	ug/l	97
46) Dibromomethane	7.568	93	9236	4.859	ug/l	95
47) Bromodichloromethane	7.799	83	19311	4.873	ug/l	98
48) Methyl methacrylate	7.671	41	14173	4.603	ug/l	95
49) 1,4-Dioxane	7.641	88	3770	89.028	ug/l #	91
51) 4-Methyl-2-Pentanone	8.555	43	97596	25.688	ug/l	97
52) Toluene	8.702	92	28194	4.662	ug/l	100
53) t-1,3-Dichloropropene	8.964	75	15951	4.231	ug/l	95
54) cis-1,3-Dichloropropene	8.348	75	19250	4.751	ug/l	97
55) 1,1,2-Trichloroethane	9.134	97	11325	4.654	ug/l	96
56) Ethyl methacrylate	9.098	69	15799	4.148	ug/l	96
57) 1,3-Dichloropropane	9.293	76	18244	4.509	ug/l	99
58) 2-Chloroethyl Vinyl ether	8.226	63	52878	25.688	ug/l	95
59) 2-Hexanone	9.415	43	55721	20.893	ug/l	97
60) Dibromochloromethane	9.506	129	13990	4.650	ug/l	99
61) 1,2-Dibromoethane	9.592	107	11968	4.545	ug/l	100
64) Tetrachloroethene	9.256	164	11027	5.035	ug/l	93
65) Chlorobenzene	10.061	112	32494	4.941	ug/l	97
66) 1,1,1,2-Tetrachloroethane	10.140	131	11291	4.800	ug/l	95
67) Ethyl Benzene	10.177	91	51672	4.815	ug/l	99
68) m/p-Xylenes	10.281	106	40552	9.701	ug/l	99
69) o-Xylene	10.622	106	19035	4.860	ug/l	98
70) Styrene	10.640	104	32097	4.773	ug/l	98
71) Bromoform	10.781	173	9949	4.738	ug/l #	99
73) Isopropylbenzene	10.945	105	50807	4.844	ug/l	99
74) N-amyl acetate	10.829	43	19243	4.223	ug/l	96
75) 1,1,2,2-Tetrachloroethane	11.195	83	16109	4.552	ug/l	96
76) 1,2,3-Trichloropropane	11.226	75	14100m	4.722	ug/l	
77) Bromobenzene	11.183	156	13538	4.836	ug/l	91
78) n-propylbenzene	11.287	91	57143	4.704	ug/l	97
79) 2-Chlorotoluene	11.348	91	35216	4.754	ug/l	98
80) 1,3,5-Trimethylbenzene	11.433	105	41155	4.774	ug/l	97
81) trans-1,4-Dichloro-2-b...	11.000	75	4948	4.055	ug/l #	87
82) 4-Chlorotoluene	11.439	91	39681	4.591	ug/l	94
83) tert-Butylbenzene	11.695	119	43520	4.858	ug/l	99
84) 1,2,4-Trimethylbenzene	11.732	105	40951	4.709	ug/l	97
85) sec-Butylbenzene	11.872	105	53557	4.933	ug/l	98
86) p-Isopropyltoluene	11.988	119	45373	4.890	ug/l	99
87) 1,3-Dichlorobenzene	11.951	146	25073	4.859	ug/l	99
88) 1,4-Dichlorobenzene	12.024	146	25523	4.786	ug/l	92
89) n-Butylbenzene	12.317	91	40625	4.676	ug/l	100
90) Hexachloroethane	12.518	117	8642	4.563	ug/l	95
91) 1,2-Dichlorobenzene	12.317	146	23201	4.637	ug/l	100
92) 1,2-Dibromo-3-Chloropr...	12.927	75	3579	4.156	ug/l	97
93) 1,2,4-Trichlorobenzene	13.573	180	14300	4.328	ug/l	97
94) Hexachlorobutadiene	13.707	225	6803	4.977	ug/l	96
95) Naphthalene	13.756	128	40122	3.182	ug/l	98
96) 1,2,3-Trichlorobenzene	13.945	180	13011	4.270	ug/l	99

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120425\
Data File : VX048715.D
Acq On : 04 Dec 2025 09:16
Operator : JC/MD
Sample : VSTDICC005
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 3 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
VSTDICC005

Manual Integrations
APPROVED

Reviewed By :John Carlone 12/05/2025
Supervised By :Semsettin Yesilyurt 12/05/2025

Quant Time: Dec 05 02:07:22 2025
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X120425W.M
Quant Title : SW846 8260
QLast Update : Fri Dec 05 02:05:18 2025
Response via : Initial Calibration

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

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

Instrument :
MSVOA_X
ClientSampleId :
VSTDICC005

Manual Integrations APPROVED

Abundance

TIC: VX048715.D\data.ms

Time-->

Chlorodifluoromethane, T
Chloromethane, T
Chloromethane, T
Trichlorofluoromethane, T
Dichloromethane, T
Acetone, T
Methyl acetate, T
Methyl acetate, T
Methyl acetate, T
Tert butyl alcohol, T
1,1-Dichloroethane, P
Diisopropyl ether, T
2,2-Dichloropropane, T
Ethyl acetate, T
Bromochloromethane, T
Chloroform, T
Dichloromethane, T
Carbon tetrachloride, T
1,2-Dichloroethane, d4, S
1,2-Dichloroethane, TM
Isopropyl Acetate, T
1,4-Difluorobenzene, I
Trichloroethene, TM
1,2-Dichlorobenzene, T
1,4-Dichlorobenzene, T
Bromodichloromethane, T
cis-1,3-Dichloropropene, T
2-Chloroethyl Vinyl ether, T
4-Methyl-2-Pentanone, T
Toluene, T
1,3-Dichloropropene, T
Ethyl acetate, T
1,3-Dichlorobenzene, T
1,2-Dichlorobenzene, T
1,2-Dichlorobenzene, T
Chlorobenzene, PM
1,1,2-Trichloroethane, T
m/p-Xylenes, T
Bromochloroacetate, T
trans-1,2-Dichloropropene, T
trans-1,2-Dichloropropene, T
1,2-Dichlorobenzene, T
1,2-Dichlorobenzene, T
1,2,4-Trichlorobenzene, T
1,3-Dichlorobenzene, T
1,3-Dichlorobenzene, T
1,2-Dichlorobenzene, T
Hexachloroethane, T
1,2-Dibromo-3-Chloropropane, T
1,2,4-Trichlorobenzene, T
1,2,4-Trichlorobenzene, I
1,2,3-Trichlorobenzene, T

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120425\
 Data File : VX048716.D
 Acq On : 04 Dec 2025 09:36
 Operator : JC/MD
 Sample : VSTDICC020
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 VSTDICC020

Manual Integrations
 APPROVED

Reviewed By : John Carlone 12/05/2025
 Supervised By : Semsettin Yesilyurt 12/05/2025

Quant Time: Dec 05 02:08:12 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X120425W.M
 Quant Title : SW846 8260
 QLast Update : Fri Dec 05 02:05:18 2025
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.538	168	219464	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.745	114	318746	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.037	117	280368	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.006	152	141102	50.000	ug/l	0.00

System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.934	65	54232	18.411	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	36.820%#
35) Dibromofluoromethane	5.373	113	45770	20.855	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	41.700%#
50) Toluene-d8	8.635	98	148035	19.243	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	38.480%#
62) 4-Bromofluorobenzene	11.061	95	53305	20.094	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	40.180%#

Target Compounds					Qvalue	
2) Dichlorodifluoromethane	1.172	85	47089	21.059	ug/l	97
3) Chloromethane	1.301	50	52390	18.567	ug/l	96
4) Vinyl Chloride	1.380	62	58519	19.473	ug/l	98
5) Bromomethane	1.618	94	37813	20.361	ug/l	99
6) Chloroethane	1.697	64	34583	18.490	ug/l	96
7) Trichlorofluoromethane	1.898	101	85315	19.808	ug/l	99
8) Diethyl Ether	2.136	74	29915	17.401	ug/l	97
9) 1,1,2-Trichlorotrifluo...	2.337	101	48761	19.529	ug/l	99
10) Methyl Iodide	2.459	142	65436	18.408	ug/l	99
11) Tert butyl alcohol	2.934	59	46577	97.940	ug/l	99
12) 1,1-Dichloroethene	2.325	96	49149	19.286	ug/l	97
13) Acrolein	2.233	56	29353	93.831	ug/l	99
14) Allyl chloride	2.666	41	85429	19.967	ug/l	97
15) Acrylonitrile	3.056	53	142975	96.485	ug/l	99
16) Acetone	2.373	43	97117	86.124	ug/l	98
17) Carbon Disulfide	2.514	76	143976	19.454	ug/l	97
18) Methyl Acetate	2.697	43	61156	19.276	ug/l	100
19) Methyl tert-butyl Ether	3.105	73	156080	18.819	ug/l	100
20) Methylene Chloride	2.788	84	51424	18.245	ug/l	97
21) trans-1,2-Dichloroethene	3.093	96	50138	19.276	ug/l	98
22) Diisopropyl ether	3.751	45	166346	20.803	ug/l	97
23) Vinyl Acetate	3.715	43	707279	105.266	ug/l	97
24) 1,1-Dichloroethane	3.605	63	91130	20.437	ug/l	99
25) 2-Butanone	4.532	43	176425	106.006	ug/l	91
26) 2,2-Dichloropropane	4.465	77	77488	19.639	ug/l	98
27) cis-1,2-Dichloroethene	4.483	96	55622	18.951	ug/l	92
28) Bromochloromethane	4.885	49	43554	19.928	ug/l	85
29) Tetrahydrofuran	4.989	42	126095	99.224	ug/l	94
30) Chloroform	5.080	83	92401	19.339	ug/l	97
31) Cyclohexane	5.458	56	84544	21.236	ug/l	95
32) 1,1,1-Trichloroethane	5.373	97	83804	19.575	ug/l	97
36) 1,1-Dichloropropene	5.678	75	65812	21.894	ug/l	97
37) Ethyl Acetate	4.696	43	83566	23.266	ug/l	97
38) Carbon Tetrachloride	5.666	117	75238	22.097	ug/l	92
39) Methylcyclohexane	7.366	83	70160	19.739	ug/l	99
40) Benzene	6.025	78	167542	18.802	ug/l	99

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120425\
 Data File : VX048716.D
 Acq On : 04 Dec 2025 09:36
 Operator : JC/MD
 Sample : VSTDICC020
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 VSTDICC020

Manual Integrations
 APPROVED

Reviewed By :John Carlone 12/05/2025
 Supervised By :Semsettin Yesilyurt 12/05/2025

Quant Time: Dec 05 02:08:12 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X120425W.M
 Quant Title : SW846 8260
 QLast Update : Fri Dec 05 02:05:18 2025
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Methacrylonitrile	4.904	41	38557	23.593	ug/l	95
42) 1,2-Dichloroethane	6.068	62	55989	17.752	ug/l	99
43) Isopropyl Acetate	6.324	43	94137	17.308	ug/l	99
44) Trichloroethene	7.110	130	46220	19.879	ug/l	99
45) 1,2-Dichloropropane	7.415	63	40049	17.982	ug/l	93
46) Dibromomethane	7.568	93	30901	18.296	ug/l	96
47) Bromodichloromethane	7.805	83	62163	17.653	ug/l	96
48) Methyl methacrylate	7.677	41	45239	16.535	ug/l	96
49) 1,4-Dioxane	7.647	88	13475	358.118	ug/l	98
51) 4-Methyl-2-Pentanone	8.555	43	299548	88.733	ug/l	100
52) Toluene	8.702	92	105894	19.707	ug/l	100
53) t-1,3-Dichloropropene	8.964	75	62519	18.662	ug/l	98
54) cis-1,3-Dichloropropene	8.348	75	67666	18.794	ug/l	93
55) 1,1,2-Trichloroethane	9.134	97	41452	19.173	ug/l	98
56) Ethyl methacrylate	9.098	69	64049	18.926	ug/l	94
57) 1,3-Dichloropropane	9.293	76	68153	18.957	ug/l	100
58) 2-Chloroethyl Vinyl ether	8.226	63	170980	93.480	ug/l	97
59) 2-Hexanone	9.415	43	220146	92.897	ug/l	98
60) Dibromochloromethane	9.506	129	52484	19.632	ug/l	100
61) 1,2-Dibromoethane	9.592	107	45333	19.373	ug/l	99
64) Tetrachloroethene	9.256	164	40200	19.671	ug/l	98
65) Chlorobenzene	10.061	112	118824	19.362	ug/l	99
66) 1,1,1,2-Tetrachloroethane	10.146	131	41197	18.768	ug/l	98
67) Ethyl Benzene	10.177	91	197228	19.695	ug/l	99
68) m/p-Xylenes	10.281	106	153388	39.322	ug/l	99
69) o-Xylene	10.622	106	73749	20.176	ug/l	98
70) Styrene	10.640	104	126646	20.182	ug/l	98
71) Bromoform	10.781	173	37588	19.181	ug/l #	97
73) Isopropylbenzene	10.945	105	192126	20.205	ug/l	99
74) N-amyl acetate	10.823	43	76117	18.423	ug/l	94
75) 1,1,2,2-Tetrachloroethane	11.195	83	58417	18.208	ug/l	97
76) 1,2,3-Trichloropropane	11.219	75	47375m	17.500	ug/l	
77) Bromobenzene	11.177	156	48752	19.209	ug/l	96
78) n-propylbenzene	11.287	91	223424	20.288	ug/l	97
79) 2-Chlorotoluene	11.347	91	132195	19.685	ug/l	98
80) 1,3,5-Trimethylbenzene	11.433	105	159778	20.443	ug/l	100
81) trans-1,4-Dichloro-2-b...	11.000	75	20567	18.590	ug/l #	78
82) 4-Chlorotoluene	11.433	91	156834	20.013	ug/l	98
83) tert-Butylbenzene	11.695	119	165236	20.346	ug/l	98
84) 1,2,4-Trimethylbenzene	11.732	105	158676	20.126	ug/l	100
85) sec-Butylbenzene	11.872	105	201582	20.481	ug/l	98
86) p-Isopropyltoluene	11.988	119	172223	20.471	ug/l	100
87) 1,3-Dichlorobenzene	11.951	146	92004	19.665	ug/l	98
88) 1,4-Dichlorobenzene	12.024	146	90959	18.813	ug/l	98
89) n-Butylbenzene	12.317	91	173673	22.048	ug/l	100
90) Hexachloroethane	12.518	117	35390	20.611	ug/l	100
91) 1,2-Dichlorobenzene	12.317	146	96083	21.183	ug/l	98
92) 1,2-Dibromo-3-Chloropr...	12.926	75	15398	19.720	ug/l	97
93) 1,2,4-Trichlorobenzene	13.567	180	58891	19.658	ug/l	98
94) Hexachlorobutadiene	13.707	225	25503	20.581	ug/l	97
95) Naphthalene	13.756	128	185802	23.652	ug/l	100
96) 1,2,3-Trichlorobenzene	13.938	180	55848	20.215	ug/l	99

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120425\
Data File : VX048716.D
Acq On : 04 Dec 2025 09:36
Operator : JC/MD
Sample : VSTDICC020
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 4 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
VSTDICC020

Manual Integrations
APPROVED

Reviewed By :John Carlone 12/05/2025
Supervised By :Semsettin Yesilyurt 12/05/2025

Quant Time: Dec 05 02:08:12 2025
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X120425W.M
Quant Title : SW846 8260
QLast Update : Fri Dec 05 02:05:18 2025
Response via : Initial Calibration

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

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120425\
 Data File : VX048716.D
 Acq On : 04 Dec 2025 09:36
 Operator : JC/MD
 Sample : VSTDIC020
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 4 Sample Multiplier: 1

Instrument :

MSVOA_X

Client SampleId :

VSTDIC020

Manual Integrations

APPROVED

Reviewed By :John Carlone 12/05/2025

Supervised By :Semsettin Yesilyurt 12/05/2025

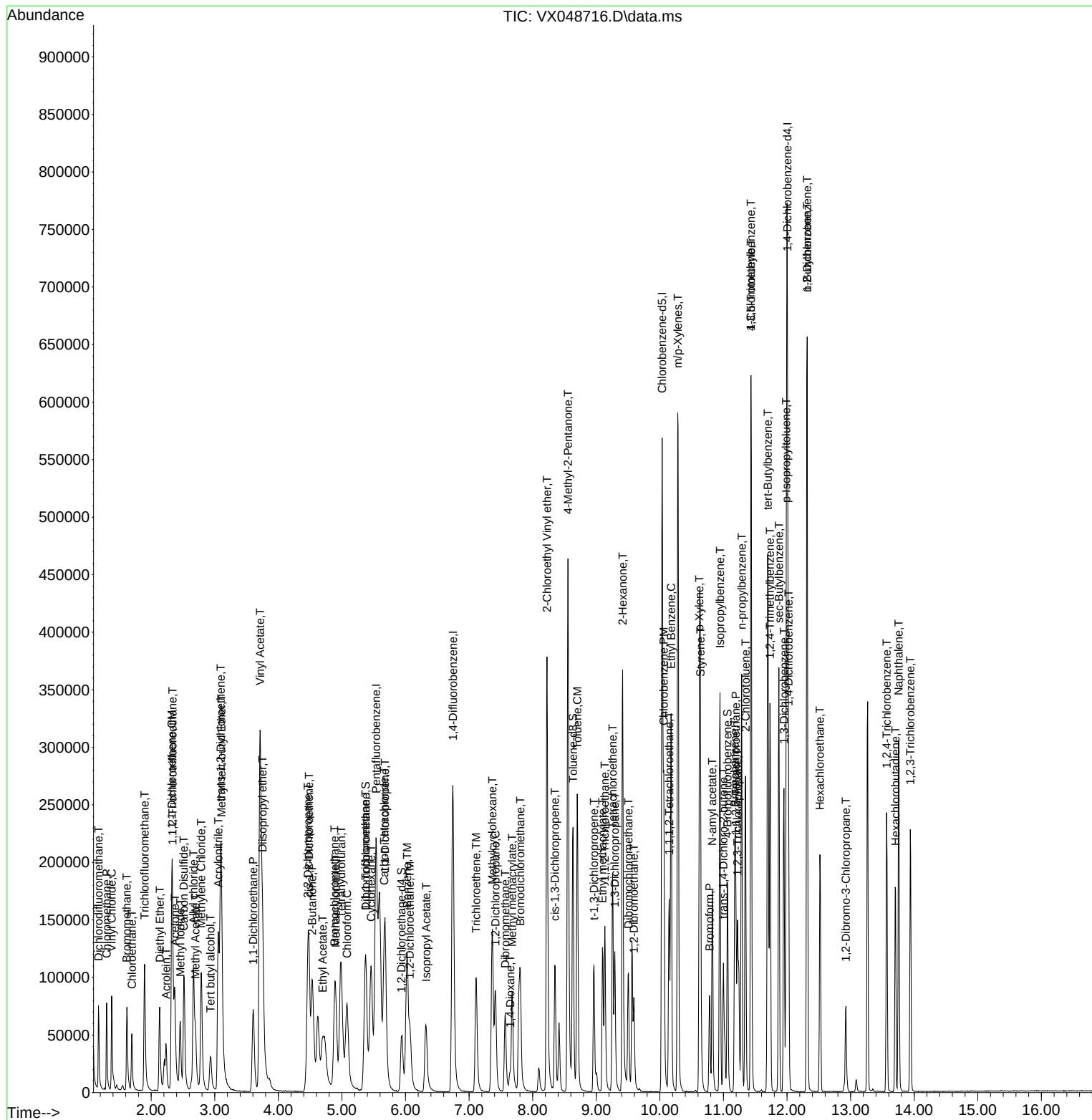
Quant Time: Dec 05 02:08:12 2025

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X120425W.M

Quant Title : SW846 8260

QLast Update : Fri Dec 05 02:05:18 2025

Response via : Initial Calibration



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120425\
 Data File : VX048717.D
 Acq On : 04 Dec 2025 09:57
 Operator : JC/MD
 Sample : VSTDICCC050
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 VSTDICCC050

Manual Integrations
 APPROVED

Reviewed By : John Carlone 12/05/2025
 Supervised By : Semsettin Yesilyurt 12/05/2025

Quant Time: Dec 05 02:09:00 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X120425W.M
 Quant Title : SW846 8260
 QLast Update : Fri Dec 05 02:05:18 2025
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.538	168	186557	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.745	114	289095	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.037	117	279030	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.006	152	136461	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.940	65	107907	43.095	ug/l	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery =	86.180%		
35) Dibromofluoromethane	5.373	113	96652	48.556	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery =	97.120%		
50) Toluene-d8	8.635	98	379431	54.380	ug/l	0.00
Spiked Amount 50.000	Range 92 - 112		Recovery =	108.760%		
62) 4-Bromofluorobenzene	11.061	95	132614	55.117	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery =	110.240%		
Target Compounds					Qvalue	
2) Dichlorodifluoromethane	1.173	85	106688	56.128	ug/l	100
3) Chloromethane	1.301	50	120299	50.154	ug/l	100
4) Vinyl Chloride	1.380	62	132300	51.790	ug/l	100
5) Bromomethane	1.618	94	87152	55.207	ug/l	100
6) Chloroethane	1.697	64	79501	50.005	ug/l	100
7) Trichlorofluoromethane	1.898	101	192516	52.582	ug/l	100
8) Diethyl Ether	2.136	74	70921	48.532	ug/l	100
9) 1,1,2-Trichlorotrifluo...	2.337	101	111195	52.388	ug/l	100
10) Methyl Iodide	2.459	142	161441	53.427	ug/l	100
11) Tert butyl alcohol	2.934	59	107184	265.136	ug/l	100
12) 1,1-Dichloroethene	2.325	96	112110	51.751	ug/l	100
13) Acrolein	2.233	56	79940	212.073	ug/l	100
14) Allyl chloride	2.666	41	200567	55.145	ug/l	100
15) Acrylonitrile	3.056	53	342616	271.992	ug/l	100
16) Acetone	2.374	43	229958	239.900	ug/l	100
17) Carbon Disulfide	2.520	76	330819	52.586	ug/l	100
18) Methyl Acetate	2.697	43	145030	53.775	ug/l	100
19) Methyl tert-butyl Ether	3.105	73	372341	52.813	ug/l	100
20) Methylene Chloride	2.788	84	121277	50.618	ug/l	100
21) trans-1,2-Dichloroethene	3.093	96	116779	52.817	ug/l	100
22) Diisopropyl ether	3.751	45	292407	43.018	ug/l	100
23) Vinyl Acetate	3.715	43	1289003	225.685	ug/l	100
24) 1,1-Dichloroethane	3.605	63	169753	44.783	ug/l	100
25) 2-Butanone	4.538	43	314742	222.473	ug/l	100
26) 2,2-Dichloropropane	4.471	77	155016	46.219	ug/l	100
27) cis-1,2-Dichloroethene	4.483	96	117259	46.998	ug/l	100
28) Bromochloromethane	4.885	49	77742	41.844	ug/l	100
29) Tetrahydrofuran	4.983	42	231381	214.189	ug/l	100
30) Chloroform	5.080	83	182965	45.048	ug/l	100
31) Cyclohexane	5.464	56	148617	43.914	ug/l	100
32) 1,1,1-Trichloroethane	5.367	97	170109	46.744	ug/l	100
36) 1,1-Dichloropropene	5.684	75	126774	46.500	ug/l	100
37) Ethyl Acetate	4.696	43	156032	47.897	ug/l	100
38) Carbon Tetrachloride	5.666	117	152584	49.410	ug/l	100
39) Methylcyclohexane	7.367	83	185682	57.597	ug/l	100
40) Benzene	6.025	78	382480	47.325	ug/l	100

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120425\
 Data File : VX048717.D
 Acq On : 04 Dec 2025 09:57
 Operator : JC/MD
 Sample : VSTDICCC050
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 VSTDICCC050

Manual Integrations
 APPROVED

Reviewed By : John Carlone 12/05/2025
 Supervised By : Semsettin Yesilyurt 12/05/2025

Quant Time: Dec 05 02:09:00 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X120425W.M
 Quant Title : SW846 8260
 QLast Update : Fri Dec 05 02:05:18 2025
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Methacrylonitrile	4.904	41	70025	47.244	ug/l	100
42) 1,2-Dichloroethane	6.074	62	130710	45.694	ug/l	100
43) Isopropyl Acetate	6.324	43	218619	44.319	ug/l	100
44) Trichloroethene	7.111	130	106003	50.267	ug/l	100
45) 1,2-Dichloropropane	7.415	63	110713	54.808	ug/l	100
46) Dibromomethane	7.568	93	83655	54.610	ug/l	100
47) Bromodichloromethane	7.806	83	169698	53.133	ug/l	100
48) Methyl methacrylate	7.677	41	134523	54.212	ug/l	100
49) 1,4-Dioxane	7.641	88	36323	1064.346	ug/l	100
51) 4-Methyl-2-Pentanone	8.555	43	874869	285.735	ug/l	100
52) Toluene	8.702	92	278645	57.175	ug/l	100
53) t-1,3-Dichloropropene	8.964	75	173511	57.105	ug/l	100
54) cis-1,3-Dichloropropene	8.348	75	186114	56.995	ug/l	100
55) 1,1,2-Trichloroethane	9.135	97	109166	55.671	ug/l	100
56) Ethyl methacrylate	9.098	69	181275	59.060	ug/l	100
57) 1,3-Dichloropropane	9.293	76	181544	55.677	ug/l	100
58) 2-Chloroethyl Vinyl ether	8.226	63	496729	299.430	ug/l	100
59) 2-Hexanone	9.415	43	639088	297.342	ug/l	100
60) Dibromochloromethane	9.506	129	135150	55.738	ug/l	100
61) 1,2-Dibromoethane	9.592	107	118684	55.922	ug/l	100
64) Tetrachloroethene	9.256	164	98419	48.390	ug/l	100
65) Chlorobenzene	10.061	112	303422	49.679	ug/l	100
66) 1,1,1,2-Tetrachloroethane	10.147	131	105994	48.520	ug/l	100
67) Ethyl Benzene	10.177	91	528248	53.003	ug/l	100
68) m/p-Xylenes	10.281	106	399998	103.033	ug/l	100
69) o-Xylene	10.622	106	189164	51.999	ug/l	100
70) Styrene	10.640	104	326481	52.277	ug/l	100
71) Bromoform	10.781	173	91650	46.993	ug/l #	100
73) Isopropylbenzene	10.945	105	495714	53.906	ug/l	100
74) N-amyl acetate	10.823	43	237184	59.361	ug/l	100
75) 1,1,2,2-Tetrachloroethane	11.195	83	160851	51.840	ug/l	100
76) 1,2,3-Trichloropropane	11.220	75	130756m	49.942	ug/l	
77) Bromobenzene	11.177	156	120915	49.264	ug/l	100
78) n-propylbenzene	11.287	91	587537	55.165	ug/l	100
79) 2-Chlorotoluene	11.348	91	343958	52.961	ug/l	100
80) 1,3,5-Trimethylbenzene	11.433	105	402294	53.222	ug/l	100
81) trans-1,4-Dichloro-2-b...	11.000	75	56706	52.997	ug/l	100
82) 4-Chlorotoluene	11.433	91	407164	53.724	ug/l	100
83) tert-Butylbenzene	11.695	119	412891	52.570	ug/l	100
84) 1,2,4-Trimethylbenzene	11.732	105	407120	53.394	ug/l	100
85) sec-Butylbenzene	11.872	105	513956	53.995	ug/l	100
86) p-Isopropyltoluene	11.988	119	433944	53.334	ug/l	100
87) 1,3-Dichlorobenzene	11.951	146	224775	49.677	ug/l	100
88) 1,4-Dichlorobenzene	12.024	146	226733	48.491	ug/l	100
89) n-Butylbenzene	12.311	91	403752	53.000	ug/l	100
90) Hexachloroethane	12.518	117	81034	48.798	ug/l	100
91) 1,2-Dichlorobenzene	12.317	146	215818	49.198	ug/l	100
92) 1,2-Dibromo-3-Chloropr...	12.920	75	36555	48.408	ug/l	100
93) 1,2,4-Trichlorobenzene	13.567	180	136083	46.970	ug/l	100
94) Hexachlorobutadiene	13.707	225	53860	44.943	ug/l	100
95) Naphthalene	13.756	128	428755	51.898	ug/l	100
96) 1,2,3-Trichlorobenzene	13.939	180	126875	47.485	ug/l	100

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120425\
Data File : VX048717.D
Acq On : 04 Dec 2025 09:57
Operator : JC/MD
Sample : VSTDICCC050
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 5 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
VSTDICCC050

Manual Integrations
APPROVED

Reviewed By :John Carlone 12/05/2025
Supervised By :Semsettin Yesilyurt 12/05/2025

Quant Time: Dec 05 02:09:00 2025
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X120425W.M
Quant Title : SW846 8260
QLast Update : Fri Dec 05 02:05:18 2025
Response via : Initial Calibration

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

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

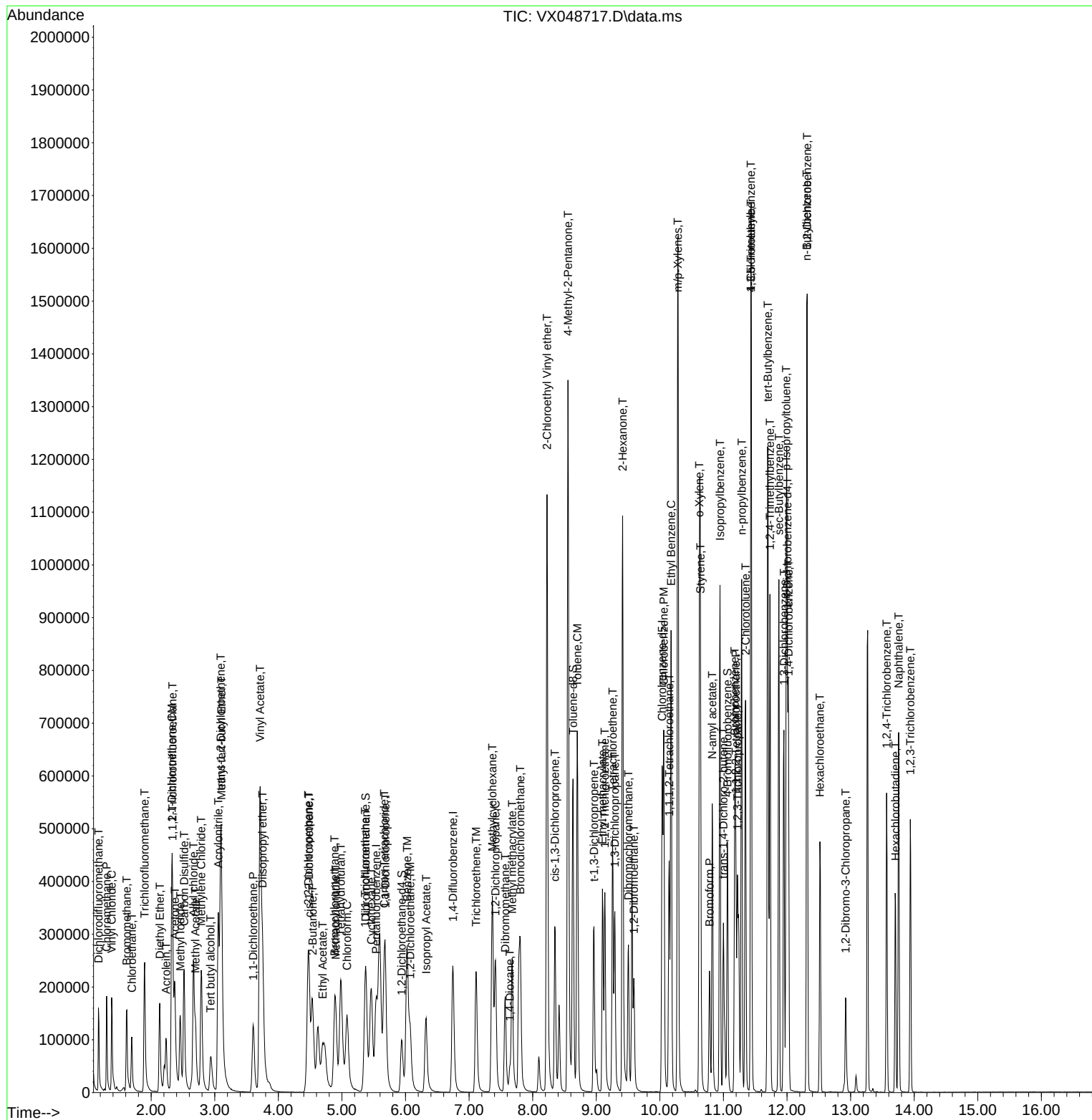
```
Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120425\
Data File : VX048717.D
Acq On    : 04 Dec 2025  09:57
Operator  : JC/MD
Sample    : VSTDICCC050
Misc      : 5.0mL/MSVOA_X/WATER
ALS Vial  : 5      Sample Multiplier: 1
```

Instrument :
MSVOA_X
ClientSampleId :
VSTDICCC050

Manual Integrations APPROVED

Reviewed By :John Carlone 12/05/2025
Supervised By :Semsettin Yesilyurt 12/05/2025

Quant Time: Dec 05 02:09:00 2025
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X120425W.M
Quant Title : SW846 8260
QLast Update : Fri Dec 05 02:05:18 2025
Response via : Initial Calibration



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120425\
 Data File : VX048718.D
 Acq On : 04 Dec 2025 10:17
 Operator : JC/MD
 Sample : VSTDICC100
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 VSTDICC100

Manual Integrations
 APPROVED

Reviewed By : John Carlone 12/05/2025
 Supervised By : Semsettin Yesilyurt 12/05/2025

Quant Time: Dec 05 02:09:49 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X120425W.M
 Quant Title : SW846 8260
 QLast Update : Fri Dec 05 02:05:18 2025
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.537	168	149943	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.745	114	269490	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.037	117	222177	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.006	152	112983	50.000	ug/l	0.00

System Monitoring Compounds

33) 1,2-Dichloroethane-d4	5.940	65	236091	117.312	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	= 234.620%#	
35) Dibromofluoromethane	5.373	113	191311	103.102	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 206.200%#	
50) Toluene-d8	8.634	98	566679	87.124	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	= 174.240%#	
62) 4-Bromofluorobenzene	11.061	95	213976	95.403	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 190.800%#	

Target Compounds

						Qvalue
2) Dichlorodifluoromethane	1.172	85	148747	97.363	ug/l	98
3) Chloromethane	1.300	50	201139	104.333	ug/l	99
4) Vinyl Chloride	1.380	62	209791	102.178	ug/l	100
5) Bromomethane	1.611	94	132130	104.136	ug/l	100
6) Chloroethane	1.691	64	130144	101.847	ug/l	98
7) Trichlorofluoromethane	1.892	101	282037	95.844	ug/l	97
8) Diethyl Ether	2.136	74	130414	111.035	ug/l	100
9) 1,1,2-Trichlorotrifluo...	2.337	101	156439	91.702	ug/l	98
10) Methyl Iodide	2.459	142	286982	118.165	ug/l	100
11) Tert butyl alcohol	2.959	59	204928	630.705	ug/l	100
12) 1,1-Dichloroethene	2.325	96	175195	100.620	ug/l	96
13) Acrolein	2.233	56	175644	510.093	ug/l	99
14) Allyl chloride	2.666	41	351406	120.211	ug/l	99
15) Acrylonitrile	3.056	53	634692	626.899	ug/l	99
16) Acetone	2.373	43	428956	556.775	ug/l	98
17) Carbon Disulfide	2.514	76	519336	102.710	ug/l	99
18) Methyl Acetate	2.696	43	271853	125.412	ug/l	100
19) Methyl tert-butyl Ether	3.105	73	675620	119.231	ug/l	99
20) Methylene Chloride	2.788	84	215266	111.785	ug/l	99
21) trans-1,2-Dichloroethene	3.093	96	192558	108.358	ug/l	98
22) Diisopropyl ether	3.751	45	683700	125.144	ug/l	98
23) Vinyl Acetate	3.715	43	3105693	676.537	ug/l	98
24) 1,1-Dichloroethane	3.605	63	366722	120.371	ug/l	98
25) 2-Butanone	4.538	43	759593	668.021	ug/l	94
26) 2,2-Dichloropropane	4.471	77	298026	110.556	ug/l	100
27) cis-1,2-Dichloroethene	4.483	96	235515	117.446	ug/l	95
28) Bromochloromethane	4.885	49	183158	122.656	ug/l	86
29) Tetrahydrofuran	4.983	42	544146	626.716	ug/l	94
30) Chloroform	5.080	83	363457	111.339	ug/l	99
31) Cyclohexane	5.464	56	258069	94.876	ug/l	94
32) 1,1,1-Trichloroethane	5.373	97	303648	103.814	ug/l	96
36) 1,1-Dichloropropene	5.684	75	227212	89.403	ug/l	99
37) Ethyl Acetate	4.696	43	370775	122.096	ug/l	99
38) Carbon Tetrachloride	5.665	117	256702	89.173	ug/l	97
39) Methylcyclohexane	7.366	83	257594	85.717	ug/l	97
40) Benzene	6.025	78	755535	100.284	ug/l	98

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120425\
 Data File : VX048718.D
 Acq On : 04 Dec 2025 10:17
 Operator : JC/MD
 Sample : VSTDICC100
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 VSTDICC100

Manual Integrations
 APPROVED

Reviewed By : John Carlone 12/05/2025
 Supervised By : Semsettin Yesilyurt 12/05/2025

Quant Time: Dec 05 02:09:49 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X120425W.M
 Quant Title : SW846 8260
 QLast Update : Fri Dec 05 02:05:18 2025
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Methacrylonitrile	4.897	41	165473	119.761	ug/l	95
42) 1,2-Dichloroethane	6.074	62	282907	106.094	ug/l	97
43) Isopropyl Acetate	6.324	43	522170	113.557	ug/l	96
44) Trichloroethene	7.110	130	186139	94.688	ug/l	94
45) 1,2-Dichloropropane	7.415	63	193808	102.923	ug/l	99
46) Dibromomethane	7.568	93	147078	102.996	ug/l	98
47) Bromodichloromethane	7.805	83	299209	100.499	ug/l	98
48) Methyl methacrylate	7.677	41	258497	111.752	ug/l	97
49) 1,4-Dioxane	7.647	88	71374	2243.567	ug/l	97
51) 4-Methyl-2-Pentanone	8.555	43	1239027	434.110	ug/l	97
52) Toluene	8.701	92	404813	89.105	ug/l	99
53) t-1,3-Dichloropropene	8.964	75	266630	94.136	ug/l	98
54) cis-1,3-Dichloropropene	8.348	75	282606	92.840	ug/l	96
55) 1,1,2-Trichloroethane	9.134	97	167036	91.380	ug/l	99
56) Ethyl methacrylate	9.098	69	280412	98.006	ug/l	95
57) 1,3-Dichloropropane	9.293	76	276473	90.960	ug/l	99
58) 2-Chloroethyl Vinyl ether	8.226	63	691132	446.925	ug/l	97
59) 2-Hexanone	9.415	43	920873	459.615	ug/l	96
60) Dibromochloromethane	9.506	129	215867	95.503	ug/l	100
61) 1,2-Dibromoethane	9.592	107	185040	93.531	ug/l	99
64) Tetrachloroethene	9.256	164	139923	86.401	ug/l	97
65) Chlorobenzene	10.061	112	463830	95.375	ug/l	99
66) 1,1,1,2-Tetrachloroethane	10.146	131	169651	97.532	ug/l	100
67) Ethyl Benzene	10.177	91	756382	95.314	ug/l	100
68) m/p-Xylenes	10.287	106	589829	190.807	ug/l	96
69) o-Xylene	10.622	106	290379	100.247	ug/l	97
70) Styrene	10.640	104	503937	101.341	ug/l	99
71) Bromoform	10.780	173	165495	106.570	ug/l	100
73) Isopropylbenzene	10.945	105	712828	93.624	ug/l	100
74) N-amyl acetate	10.829	43	336967	101.859	ug/l	95
75) 1,1,2,2-Tetrachloroethane	11.195	83	250299	97.430	ug/l	99
76) 1,2,3-Trichloropropane	11.219	75	200074m	92.297	ug/l	
77) Bromobenzene	11.183	156	204596	100.679	ug/l	92
78) n-propylbenzene	11.286	91	809506	91.800	ug/l	98
79) 2-Chlorotoluene	11.347	91	510695	94.975	ug/l	98
80) 1,3,5-Trimethylbenzene	11.433	105	595416	95.139	ug/l	100
81) trans-1,4-Dichloro-2-b...	11.000	75	93932	106.031	ug/l	90
82) 4-Chlorotoluene	11.433	91	602673	96.045	ug/l	98
83) tert-Butylbenzene	11.695	119	619946	95.335	ug/l	96
84) 1,2,4-Trimethylbenzene	11.731	105	613436	97.170	ug/l	99
85) sec-Butylbenzene	11.872	105	714356	90.644	ug/l	98
86) p-Isopropyltoluene	11.987	119	629447	93.439	ug/l	99
87) 1,3-Dichlorobenzene	11.951	146	364260	97.233	ug/l	99
88) 1,4-Dichlorobenzene	12.024	146	370371	95.671	ug/l	99
89) n-Butylbenzene	12.311	91	555297	88.041	ug/l	99
90) Hexachloroethane	12.518	117	121649	88.478	ug/l	94
91) 1,2-Dichlorobenzene	12.317	146	351927	96.897	ug/l	100
92) 1,2-Dibromo-3-Chloropr...	12.920	75	59671	95.439	ug/l	90
93) 1,2,4-Trichlorobenzene	13.566	180	240057	100.075	ug/l	99
94) Hexachlorobutadiene	13.707	225	88160	88.851	ug/l	96
95) Naphthalene	13.755	128	790203	97.658	ug/l	100
96) 1,2,3-Trichlorobenzene	13.938	180	226988	102.608	ug/l	99

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120425\
Data File : VX048718.D
Acq On : 04 Dec 2025 10:17
Operator : JC/MD
Sample : VSTDICC100
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
VSTDICC100

Manual Integrations
APPROVED

Reviewed By :John Carlone 12/05/2025
Supervised By :Semsettin Yesilyurt 12/05/2025

Quant Time: Dec 05 02:09:49 2025
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X120425W.M
Quant Title : SW846 8260
QLast Update : Fri Dec 05 02:05:18 2025
Response via : Initial Calibration

Compound	R.T.	Q	Ion	Response	Conc	Units	Dev(Min)
----------	------	---	-----	----------	------	-------	----------

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120425\
 Data File : VX048718.D
 Acq On : 04 Dec 2025 10:17
 Operator : JC/MD
 Sample : VSTDIC100
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 6 Sample Multiplier: 1

Instrument :

MSVOA_X

Client SampleId :

VSTDIC100

Manual Integrations

APPROVED

Reviewed By :John Carlone 12/05/2025

Supervised By :Semsettin Yesilyurt 12/05/2025

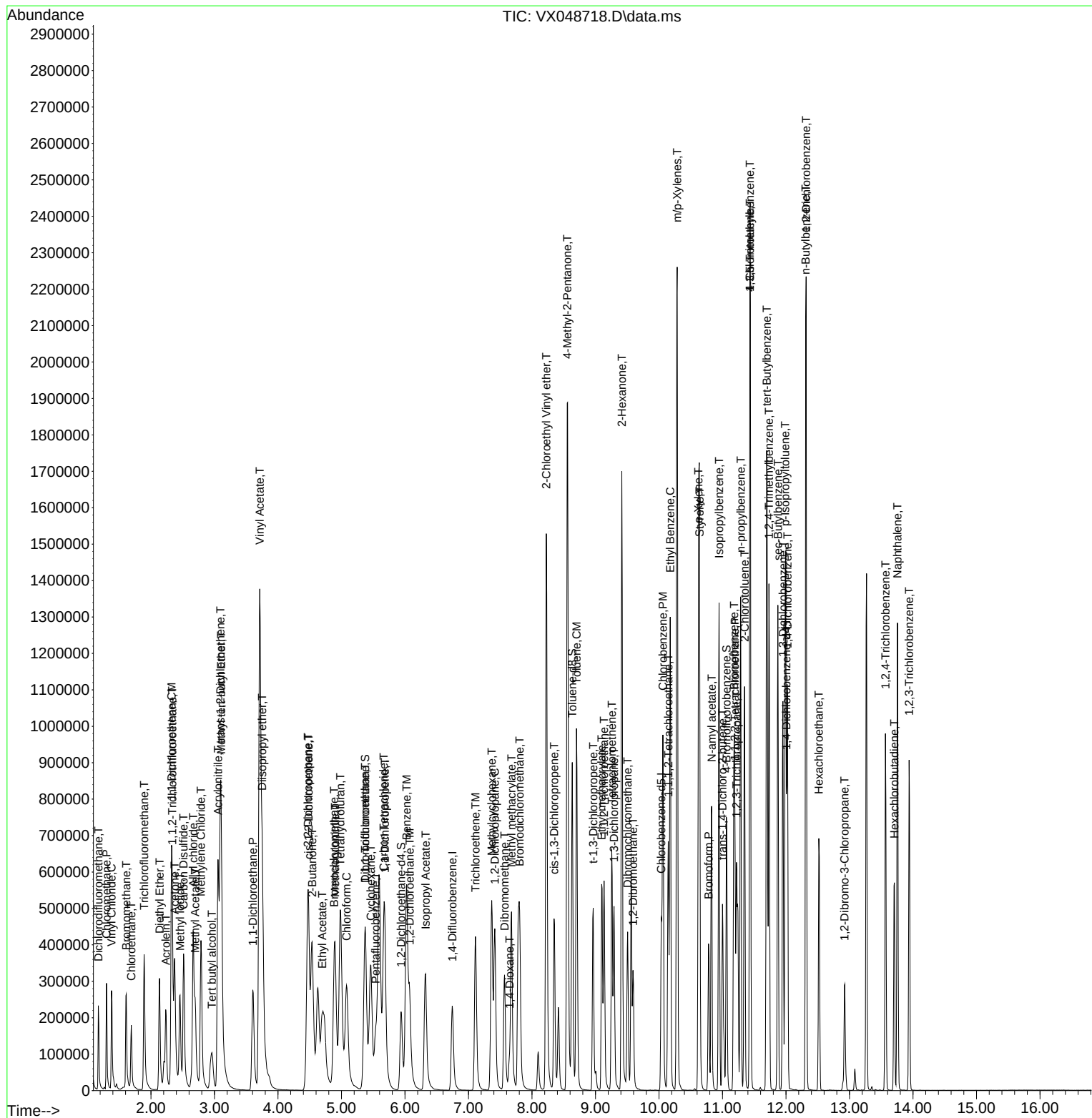
Quant Time: Dec 05 02:09:49 2025

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X120425W.M

Quant Title : SW846 8260

QLast Update : Fri Dec 05 02:05:18 2025

Response via : Initial Calibration



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120425\
 Data File : VX048719.D
 Acq On : 04 Dec 2025 10:38
 Operator : JC/MD
 Sample : VSTDICC150
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 VSTDICC150

Manual Integrations
 APPROVED

Reviewed By : John Carlone 12/05/2025
 Supervised By : Semsettin Yesilyurt 12/05/2025

Quant Time: Dec 05 02:10:38 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X120425W.M
 Quant Title : SW846 8260
 QLast Update : Fri Dec 05 02:05:18 2025
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.544	168	161225	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.745	114	266084	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.037	117	202872	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.006	152	103615	50.000	ug/l	# 0.00

System Monitoring Compounds

33) 1,2-Dichloroethane-d4	5.940	65	336173	155.353	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	= 310.700%#	
35) Dibromofluoromethane	5.373	113	284768	155.433	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 310.860%#	
50) Toluene-d8	8.635	98	980276	152.642	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	= 305.280%#	
62) 4-Bromofluorobenzene	11.061	95	311931	140.857	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 281.720%#	

Target Compounds

						Qvalue
2) Dichlorodifluoromethane	1.173	85	260977	158.870	ug/l	99
3) Chloromethane	1.301	50	321280	154.990	ug/l	99
4) Vinyl Chloride	1.380	62	340742	154.345	ug/l	99
5) Bromomethane	1.605	94	163576	119.898	ug/l	100
6) Chloroethane	1.685	64	203254	147.930	ug/l	98
7) Trichlorofluoromethane	1.892	101	488765	154.472	ug/l	97
8) Diethyl Ether	2.136	74	189537	150.080	ug/l	98
9) 1,1,2-Trichlorotrifluo...	2.337	101	282907	154.231	ug/l	99
10) Methyl Iodide	2.453	142	431805	165.355	ug/l	100
11) Tert butyl alcohol	2.977	59	236315	676.410	ug/l	100
12) 1,1-Dichloroethene	2.325	96	289225	154.488	ug/l	97
13) Acrolein	2.233	56	287608	755.796	ug/l	99
14) Allyl chloride	2.666	41	386590	122.993	ug/l	96
15) Acrylonitrile	3.056	53	695550	638.935	ug/l	99
16) Acetone	2.373	43	617325	745.203	ug/l	99
17) Carbon Disulfide	2.514	76	813894	149.701	ug/l	99
18) Methyl Acetate	2.697	43	290410	124.598	ug/l	96
19) Methyl tert-butyl Ether	3.111	73	820601	134.683	ug/l	97
20) Methylene Chloride	2.788	84	258488	124.837	ug/l	93
21) trans-1,2-Dichloroethene	3.093	96	253511	132.675	ug/l	93
22) Diisopropyl ether	3.751	45	760942	129.535	ug/l	98
23) Vinyl Acetate	3.715	43	3401948	689.215	ug/l	100
24) 1,1-Dichloroethane	3.605	63	441859	134.884	ug/l	99
25) 2-Butanone	4.544	43	831570	680.145	ug/l	100
26) 2,2-Dichloropropane	4.471	77	405592	139.930	ug/l	99
27) cis-1,2-Dichloroethene	4.477	96	304797	141.360	ug/l	100
28) Bromochloromethane	4.891	49	233635	145.511	ug/l	95
29) Tetrahydrofuran	4.983	42	740891	793.604	ug/l	96
30) Chloroform	5.080	83	542284	154.495	ug/l	99
31) Cyclohexane	5.464	56	457324	156.365	ug/l	98
32) 1,1,1-Trichloroethane	5.373	97	494982	157.387	ug/l	98
36) 1,1-Dichloropropene	5.684	75	374749	149.343	ug/l	99
37) Ethyl Acetate	4.696	43	374331	124.844	ug/l	97
38) Carbon Tetrachloride	5.666	117	429152	150.987	ug/l	97
39) Methylcyclohexane	7.367	83	475877	160.380	ug/l	99
40) Benzene	6.025	78	1152810	154.974	ug/l	99

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120425\
 Data File : VX048719.D
 Acq On : 04 Dec 2025 10:38
 Operator : JC/MD
 Sample : VSTDICC150
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 VSTDICC150

Quant Time: Dec 05 02:10:38 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X120425W.M
 Quant Title : SW846 8260
 QLast Update : Fri Dec 05 02:05:18 2025
 Response via : Initial Calibration

Manual Integrations
 APPROVED

Reviewed By :John Carlone 12/05/2025
 Supervised By :Semsettin Yesilyurt 12/05/2025

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Methacrylonitrile	4.910	41	210934	154.617	ug/l	99
42) 1,2-Dichloroethane	6.074	62	400347	152.058	ug/l	98
43) Isopropyl Acetate	6.324	43	726750	160.070	ug/l	99
44) Trichloroethene	7.110	130	308803	159.098	ug/l	99
45) 1,2-Dichloropropane	7.415	63	287426	154.593	ug/l	99
46) Dibromomethane	7.568	93	216552	153.589	ug/l	98
47) Bromodichloromethane	7.805	83	440047	149.695	ug/l	97
48) Methyl methacrylate	7.677	41	361697	158.368	ug/l	99
49) 1,4-Dioxane	7.653	88	102268	3255.837	ug/l	96
51) 4-Methyl-2-Pentanone	8.561	43	2202246	781.463	ug/l	99
52) Toluene	8.702	92	709062	158.073	ug/l	100
53) t-1,3-Dichloropropene	8.964	75	463456	165.721	ug/l	100
54) cis-1,3-Dichloropropene	8.354	75	489309	162.802	ug/l	96
55) 1,1,2-Trichloroethane	9.134	97	277366	153.680	ug/l	98
56) Ethyl methacrylate	9.104	69	481264	170.358	ug/l	97
57) 1,3-Dichloropropane	9.293	76	462553	154.128	ug/l	100
58) 2-Chloroethyl Vinyl ether	8.232	63	1324521	867.473	ug/l	99
59) 2-Hexanone	9.415	43	1628603	823.253	ug/l	99
60) Dibromochloromethane	9.506	129	344385	154.312	ug/l	99
61) 1,2-Dibromoethane	9.592	107	296681	151.881	ug/l	99
64) Tetrachloroethene	9.256	164	249690	168.852	ug/l	99
65) Chlorobenzene	10.061	112	677420	152.550	ug/l	98
66) 1,1,1,2-Tetrachloroethane	10.147	131	249387	157.014	ug/l	99
67) Ethyl Benzene	10.177	91	1156627	159.619	ug/l	99
68) m/p-Xylenes	10.287	106	895563	317.279	ug/l	97
69) o-Xylene	10.622	106	433369	163.848	ug/l	97
70) Styrene	10.640	104	744510	163.967	ug/l	99
71) Bromoform	10.781	173	240629	169.697	ug/l	100
73) Isopropylbenzene	10.945	105	1114762	159.651	ug/l	100
74) N-amyl acetate	10.829	43	488735	161.092	ug/l	95
75) 1,1,2,2-Tetrachloroethane	11.195	83	363135	154.132	ug/l	99
76) 1,2,3-Trichloropropane	11.226	75	281629m	141.666	ug/l	
77) Bromobenzene	11.183	156	299915	160.928	ug/l	95
78) n-propylbenzene	11.287	91	1284705	158.861	ug/l	98
79) 2-Chlorotoluene	11.348	91	774892	157.137	ug/l	98
80) 1,3,5-Trimethylbenzene	11.433	105	925692	161.286	ug/l	100
81) trans-1,4-Dichloro-2-b...	11.000	75	138844	170.898	ug/l	86
82) 4-Chlorotoluene	11.439	91	912659	158.596	ug/l	98
83) tert-Butylbenzene	11.695	119	977059	163.836	ug/l	97
84) 1,2,4-Trimethylbenzene	11.732	105	944577	163.151	ug/l	99
85) sec-Butylbenzene	11.872	105	1158976	160.357	ug/l	98
86) p-Isopropyltoluene	11.994	119	1005907	162.823	ug/l	99
87) 1,3-Dichlorobenzene	11.951	146	545971	158.914	ug/l	100
88) 1,4-Dichlorobenzene	12.024	146	551782	155.418	ug/l	98
89) n-Butylbenzene	12.317	91	904369	156.349	ug/l	99
90) Hexachloroethane	12.518	117	223207	177.022	ug/l	100
91) 1,2-Dichlorobenzene	12.317	146	522301	156.808	ug/l	99
92) 1,2-Dibromo-3-Chloropr...	12.920	75	103188	179.963	ug/l	96
93) 1,2,4-Trichlorobenzene	13.567	180	402548	182.988	ug/l	98
94) Hexachlorobutadiene	13.707	225	143127	157.291	ug/l	98
95) Naphthalene	13.756	128	1330400	150.662	ug/l	100
96) 1,2,3-Trichlorobenzene	13.938	180	373809	184.254	ug/l	99

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120425\
Data File : VX048719.D
Acq On : 04 Dec 2025 10:38
Operator : JC/MD
Sample : VSTDICC150
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
VSTDICC150

Manual Integrations
APPROVED

Reviewed By :John Carlone 12/05/2025
Supervised By :Semsettin Yesilyurt 12/05/2025

Quant Time: Dec 05 02:10:38 2025
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X120425W.M
Quant Title : SW846 8260
QLast Update : Fri Dec 05 02:05:18 2025
Response via : Initial Calibration

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

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

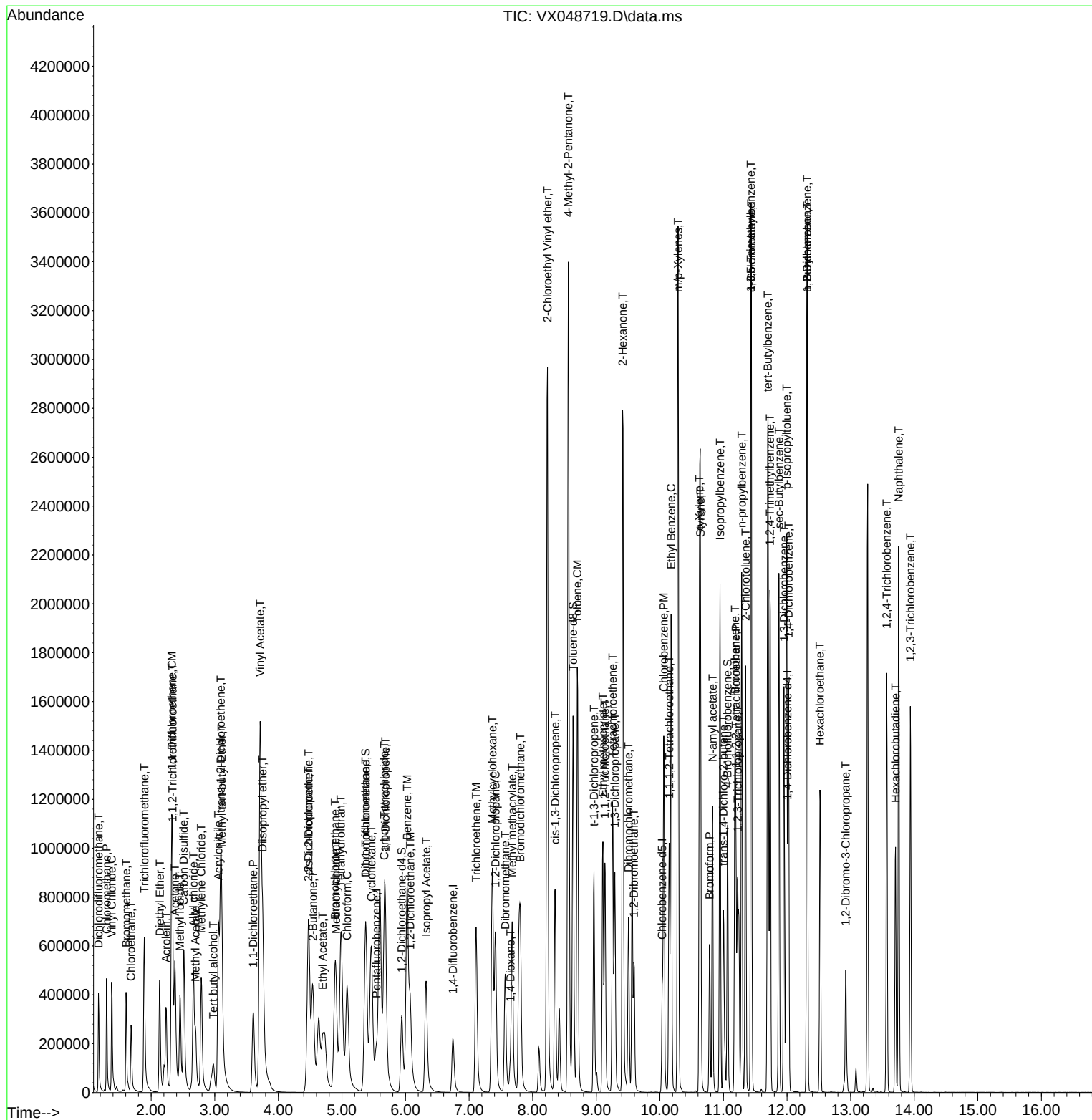
Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120425\
 Data File : VX048719.D
 Acq On : 04 Dec 2025 10:38
 Operator : JC/MD
 Sample : VSTDIC150
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 VSTDIC150

Manual Integrations APPROVED

Reviewed By : John Carlone 12/05/2025
 Supervised By : Semsettin Yesilyurt 12/05/2025

Quant Time: Dec 05 02:10:38 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X120425W.M
 Quant Title : SW846 8260
 QLast Update : Fri Dec 05 02:05:18 2025
 Response via : Initial Calibration



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120425\
 Data File : VX048721.D
 Acq On : 04 Dec 2025 14:04
 Operator : JC/MD
 Sample : VSTDICV050
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 ICVVX120425

Quant Time: Dec 05 03:31:44 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X120425W.M
 Quant Title : SW846 8260
 QLast Update : Fri Dec 05 03:27:34 2025
 Response via : Initial Calibration

Manual Integrations APPROVED

Reviewed By :John Carlone 12/05/2025
 Supervised By :Semsettin Yesilyurt 12/05/2025

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

Internal Standards						
1) Pentafluorobenzene	5.537	168	184967	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.745	114	310263	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.037	117	253965	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.000	152	132465	50.000	ug/l	0.00

System Monitoring Compounds

33) 1,2-Dichloroethane-d4	5.946	65	134442	54.154	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	= 108.300%	
35) Dibromofluoromethane	5.373	113	106728	49.960	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 99.920%	
50) Toluene-d8	8.634	98	333564	44.544	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	= 89.080%#	
62) 4-Bromofluorobenzene	11.061	95	128866	49.905	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 99.820%	

Target Compounds

						Qvalue
2) Dichlorodifluoromethane	1.172	85	110698	58.738	ug/l	98
3) Chloromethane	1.300	50	116116	48.826	ug/l	99
4) Vinyl Chloride	1.380	62	122990	48.559	ug/l	99
5) Bromomethane	1.611	94	85918	54.893	ug/l	100
6) Chloroethane	1.697	64	76727	48.675	ug/l	98
7) Trichlorofluoromethane	1.898	101	177151	48.801	ug/l	96
8) Diethyl Ether	2.136	74	69765	48.151	ug/l	100
9) 1,1,2-Trichlorotrifluo...	2.337	101	96207	45.717	ug/l	96
10) Methyl Iodide	2.459	142	150980	50.395	ug/l	99
11) Tert butyl alcohol	2.940	59	100301	250.243	ug/l	99
12) 1,1-Dichloroethene	2.325	96	100196	46.649	ug/l	93
13) Acrolein	2.233	56	88451	232.010	ug/l	99
14) Allyl chloride	2.666	41	153518	42.572	ug/l	98
15) Acrylonitrile	3.056	53	283115	226.688	ug/l	98
16) Acetone	2.373	43	194390	204.537	ug/l	98
17) Carbon Disulfide	2.520	76	256217	41.077	ug/l	100
18) Methyl Acetate	2.696	43	118728	44.401	ug/l	99
19) Methyl tert-butyl Ether	3.105	73	318949	45.629	ug/l	98
20) Methylene Chloride	2.788	84	113128	47.622	ug/l	97
21) trans-1,2-Dichloroethene	3.093	96	97130	44.308	ug/l	95
22) Diisopropyl ether	3.751	45	340522	50.527	ug/l	99
23) Vinyl Acetate	3.715	43	1563197	276.044	ug/l	98
24) 1,1-Dichloroethane	3.605	63	198478	52.811	ug/l	99
25) 2-Butanone	4.538	43	342340	244.061	ug/l	99
26) 2,2-Dichloropropane	4.471	77	152491	45.857	ug/l	99
27) cis-1,2-Dichloroethene	4.477	96	118820	48.033	ug/l	99
28) Bromochloromethane	4.885	49	82967	45.040	ug/l	97
29) Tetrahydrofuran	4.983	42	297621	277.876	ug/l	95
30) Chloroform	5.086	83	209644	52.061	ug/l	99
31) Cyclohexane	5.464	56	164989	49.171	ug/l	95
32) 1,1,1-Trichloroethane	5.373	97	180664	50.071	ug/l	97
36) 1,1-Dichloropropene	5.684	75	138860	47.458	ug/l	99
37) Ethyl Acetate	4.696	43	161815	46.283	ug/l	98
38) Carbon Tetrachloride	5.665	117	156330	47.169	ug/l	98
39) Methylcyclohexane	7.366	83	146837	42.440	ug/l	99
40) Benzene	6.025	78	436935	50.374	ug/l	99

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120425\
 Data File : VX048721.D
 Acq On : 04 Dec 2025 14:04
 Operator : JC/MD
 Sample : VSTDICV050
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 ICVVX120425

Manual Integrations
 APPROVED

Reviewed By :John Carlone 12/05/2025
 Supervised By :Semsettin Yesilyurt 12/05/2025

Quant Time: Dec 05 03:31:44 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X120425W.M
 Quant Title : SW846 8260
 QLast Update : Fri Dec 05 03:27:34 2025
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Methacrylonitrile	4.903	41	77315	48.603	ug/l	98
42) 1,2-Dichloroethane	6.074	62	156709	51.045	ug/l	97
43) Isopropyl Acetate	6.324	43	284443	53.729	ug/l	97
44) Trichloroethene	7.110	130	102412	45.250	ug/l	94
45) 1,2-Dichloropropane	7.415	63	93080	42.935	ug/l	99
46) Dibromomethane	7.568	93	71939	43.757	ug/l	95
47) Bromodichloromethane	7.805	83	147055	42.902	ug/l	99
48) Methyl methacrylate	7.677	41	112824	42.366	ug/l	96
49) 1,4-Dioxane	7.647	88	36105	985.778	ug/l	93
51) 4-Methyl-2-Pentanone	8.555	43	724550	220.496	ug/l	97
52) Toluene	8.702	92	241004	46.077	ug/l	100
53) t-1,3-Dichloropropene	8.964	75	151680	46.514	ug/l	99
54) cis-1,3-Dichloropropene	8.354	75	160155	45.699	ug/l	93
55) 1,1,2-Trichloroethane	9.134	97	96409	45.811	ug/l	99
56) Ethyl methacrylate	9.098	69	158928	48.247	ug/l	94
57) 1,3-Dichloropropane	9.293	76	158926	45.415	ug/l	100
58) 2-Chloroethyl Vinyl ether	8.226	63	431918	242.598	ug/l	98
59) 2-Hexanone	9.415	43	548159	237.637	ug/l	97
60) Dibromochloromethane	9.506	129	123372	47.409	ug/l	100
61) 1,2-Dibromoethane	9.592	107	104948	46.076	ug/l	99
64) Tetrachloroethene	9.256	164	84359	45.571	ug/l	98
65) Chlorobenzene	10.061	112	268428	48.287	ug/l	97
66) 1,1,1,2-Tetrachloroethane	10.146	131	98651	49.615	ug/l	97
67) Ethyl Benzene	10.177	91	454183	50.069	ug/l	99
68) m/p-Xylenes	10.287	106	364130	103.051	ug/l	98
69) o-Xylene	10.622	106	170785	51.580	ug/l	97
70) Styrene	10.640	104	292377	51.437	ug/l	99
71) Bromoform	10.780	173	91220	51.388	ug/l #	99
73) Isopropylbenzene	10.945	105	477424	53.483	ug/l	99
74) N-amyl acetate	10.823	43	228431	58.895	ug/l	98
75) 1,1,2,2-Tetrachloroethane	11.195	83	160485	53.282	ug/l	100
76) 1,2,3-Trichloropropane	11.219	75	131988m	53.497	ug/l	
77) Bromobenzene	11.177	156	119980	50.357	ug/l	98
78) n-propylbenzene	11.286	91	542000	52.425	ug/l	100
79) 2-Chlorotoluene	11.347	91	324683	51.501	ug/l	99
80) 1,3,5-Trimethylbenzene	11.433	105	380771	51.894	ug/l	99
81) trans-1,4-Dichloro-2-b...	11.000	75	55921	53.840	ug/l	96
82) 4-Chlorotoluene	11.433	91	386271	52.505	ug/l	100
83) tert-Butylbenzene	11.695	119	382261	50.138	ug/l	99
84) 1,2,4-Trimethylbenzene	11.731	105	381863	51.592	ug/l	99
85) sec-Butylbenzene	11.872	105	467593	50.606	ug/l	99
86) p-Isopropyltoluene	11.988	119	397475	50.326	ug/l	99
87) 1,3-Dichlorobenzene	11.951	146	215432	49.048	ug/l	99
88) 1,4-Dichlorobenzene	12.024	146	216879	47.783	ug/l	99
89) n-Butylbenzene	12.311	91	375250	50.745	ug/l	99
90) Hexachloroethane	12.518	117	75643	46.926	ug/l	97
91) 1,2-Dichlorobenzene	12.317	146	204807	48.097	ug/l	100
92) 1,2-Dibromo-3-Chloropr...	12.920	75	37176	50.715	ug/l	95
93) 1,2,4-Trichlorobenzene	13.567	180	132170	46.996	ug/l	99
94) Hexachlorobutadiene	13.707	225	52466	45.101	ug/l	95
95) Naphthalene	13.755	128	452299	55.628	ug/l	100
96) 1,2,3-Trichlorobenzene	13.938	180	128090	49.386	ug/l	98

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120425\
Data File : VX048721.D
Acq On : 04 Dec 2025 14:04
Operator : JC/MD
Sample : VSTDICV050
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
ICVVX120425

Manual Integrations
APPROVED

Reviewed By :John Carlone 12/05/2025
Supervised By :Semsettin Yesilyurt 12/05/2025

Quant Time: Dec 05 03:31:44 2025
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X120425W.M
Quant Title : SW846 8260
QLast Update : Fri Dec 05 03:27:34 2025
Response via : Initial Calibration

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

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120425\
 Data File : VX048721.D
 Acq On : 04 Dec 2025 14:04
 Operator : JC/MD
 Sample : VSTDICV050
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 13 Sample Multiplier: 1

Instrument :

MSVOA_X

Client SampleId :

ICVVX120425

Manual Integrations

APPROVED

Reviewed By : John Carlone 12/05/2025

Supervised By : Semsettin Yesilyurt 12/05/2025

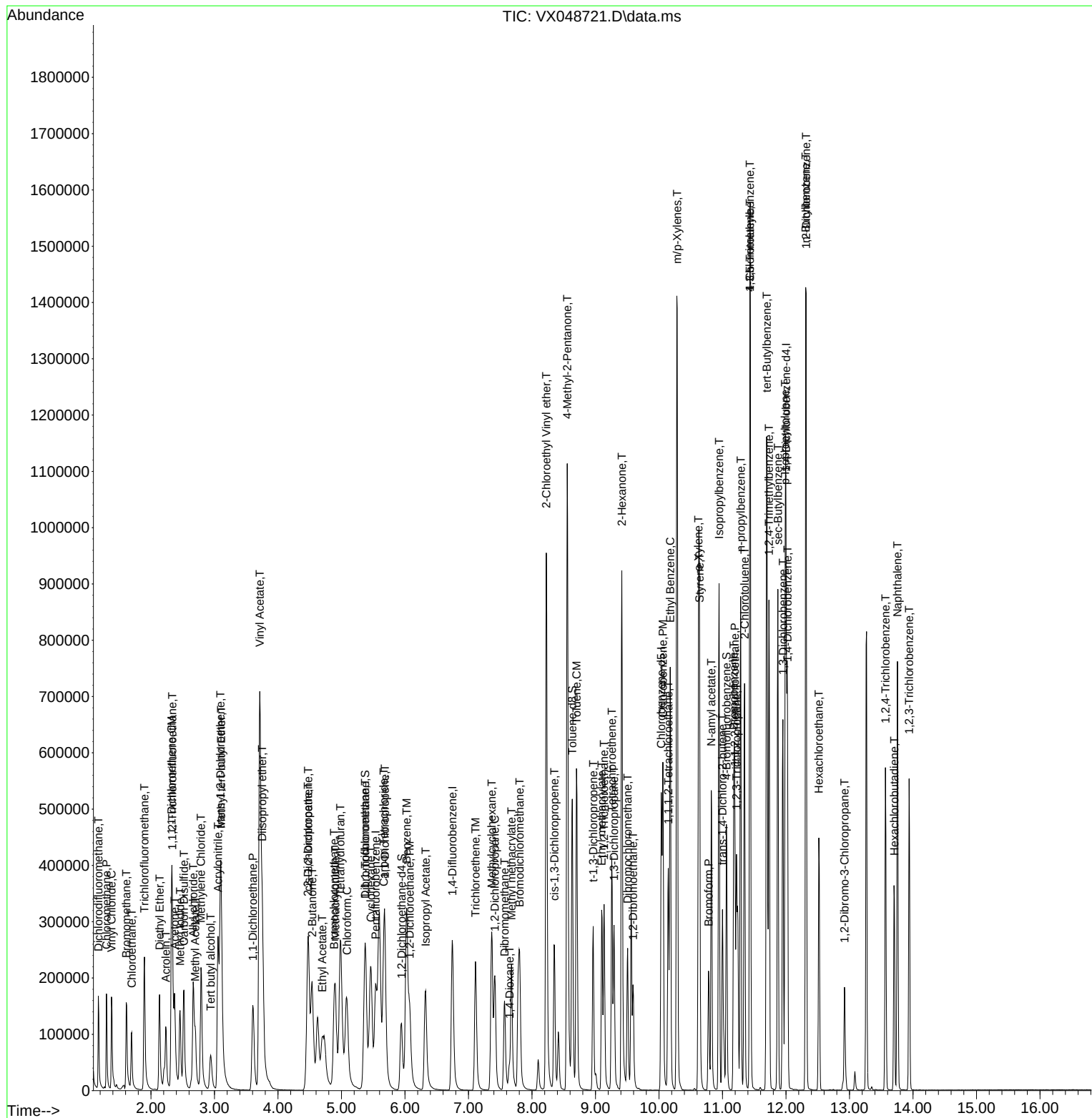
Quant Time: Dec 05 03:31:44 2025

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X120425W.M

Quant Title : SW846 8260

QLast Update : Fri Dec 05 03:27:34 2025

Response via : Initial Calibration



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120425\
 Data File : VX048721.D
 Acq On : 04 Dec 2025 14:04
 Operator : JC/MD
 Sample : VSTDICV050
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 ICVVX120425

Quant Time: Dec 05 03:31:44 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X120425W.M
 Quant Title : SW846 8260
 QLast Update : Fri Dec 05 03:27:34 2025
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
 Max. RRF Dev : 25% Max. Rel. Area : 150%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
1 I	Pentafluorobenzene	1.000	1.000	0.0	99	0.00
2 T	Dichlorodifluoromethane	0.509	0.598	-17.5	104	0.00
3 P	Chloromethane	0.643	0.628	2.3	97	0.00
4 C	Vinyl Chloride	0.685	0.665	2.9#	93	0.00
5 T	Bromomethane	0.423	0.465	-9.9	99	0.00
6 T	Chloroethane	0.426	0.415	2.6	97	0.00
7 T	Trichlorofluoromethane	0.981	0.958	2.3	92	0.00
8 T	Diethyl Ether	0.392	0.377	3.8	98	0.00
9 T	1,1,2-Trichlorotrifluoroeth	0.569	0.520	8.6	87	0.00
10 T	Methyl Iodide	0.810	0.816	-0.7	94	0.00
11 T	Tert butyl alcohol	0.108	0.108	0.0	94	0.00
12 CM	1,1-Dichloroethene	0.581	0.542	6.7#	89	0.00
13 T	Acrolein	0.091	0.096	-5.5	111	0.00
14 T	Allyl chloride	0.975	0.830	14.9	77	0.00
15 T	Acrylonitrile	0.338	0.306	9.5	83	0.00
16 T	Acetone	0.257	0.210	18.3	85	0.00
17 T	Carbon Disulfide	1.686	1.385	17.9	77	0.00
18 T	Methyl Acetate	0.723	0.642	11.2	82	0.00
19 T	Methyl tert-butyl Ether	1.890	1.724	8.8	86	0.00
20 T	Methylene Chloride	0.642	0.612	4.7	93	0.00
21 T	trans-1,2-Dichloroethene	0.593	0.525	11.5	83	0.00
22 T	Diisopropyl ether	1.822	1.841	-1.0	116	0.00
23 T	Vinyl Acetate	1.531	1.690	-10.4	121	0.00
24 P	1,1-Dichloroethane	1.016	1.073	-5.6	117	0.00
25 T	2-Butanone	0.379	0.370	2.4	109	0.00
26 T	2,2-Dichloropropane	0.899	0.824	8.3	98	0.00
27 T	cis-1,2-Dichloroethene	0.669	0.642	4.0	101	0.00
28 T	Bromochloromethane	0.498	0.449	9.8	107	0.00
29 T	Tetrahydrofuran	0.290	0.322	-11.0	129	0.00
30 C	Chloroform	1.089	1.133	-4.0#	115	0.00
31 T	Cyclohexane	0.907	0.892	1.7	111	0.00
32 T	1,1,1-Trichloroethane	0.975	0.977	-0.2	106	0.00
33 S	1,2-Dichloroethane-d4	0.671	0.727	-8.3	125	0.00
34 I	1,4-Difluorobenzene	1.000	1.000	0.0	107	0.00
35 S	Dibromofluoromethane	0.344	0.344	0.0	110	0.00
36 T	1,1-Dichloropropene	0.472	0.448	5.1	110	0.00
37 T	Ethyl Acetate	0.563	0.522	7.3	104	0.00
38 T	Carbon Tetrachloride	0.534	0.504	5.6	102	0.00
39 T	Methylcyclohexane	0.558	0.473	15.2	79	0.00
40 TM	Benzene	1.398	1.408	-0.7	114	0.00
41 T	Methacrylonitrile	0.256	0.249	2.7	110	0.00
42 TM	1,2-Dichloroethane	0.495	0.505	-2.0	120	0.00
43 T	Isopropyl Acetate	0.853	0.917	-7.5	130	0.00
44 TM	Trichloroethene	0.365	0.330	9.6	97	0.00
45 C	1,2-Dichloropropane	0.349	0.300	14.0#	84	0.00
46 T	Dibromomethane	0.265	0.232	12.5	86	0.00
47 T	Bromodichloromethane	0.552	0.474	14.1	87	0.00
48 T	Methyl methacrylate	0.429	0.364	15.2	84	0.00

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120425\
 Data File : VX048721.D
 Acq On : 04 Dec 2025 14:04
 Operator : JC/MD
 Sample : VSTDICV050
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 ICSVVX120425

Quant Time: Dec 05 03:31:44 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X120425W.M
 Quant Title : SW846 8260
 QLast Update : Fri Dec 05 03:27:34 2025
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
 Max. RRF Dev : 25% Max. Rel. Area : 150%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
49 T	1,4-Dioxane	0.006	0.006	0.0	99	0.00
50 S	Toluene-d8	1.207	1.075	10.9	88	0.00
51 T	4-Methyl-2-Pentanone	0.530	0.467	11.9	83	0.00
52 CM	Toluene	0.843	0.777	7.8#	86	0.00
53 T	t-1,3-Dichloropropene	0.526	0.489	7.0	87	0.00
54 T	cis-1,3-Dichloropropene	0.565	0.516	8.7	86	0.00
55 T	1,1,2-Trichloroethane	0.339	0.311	8.3	88	0.00
56 T	Ethyl methacrylate	0.531	0.512	3.6	88	0.00
57 T	1,3-Dichloropropane	0.564	0.512	9.2	88	0.00
58 T	2-Chloroethyl Vinyl ether	0.287	0.278	3.1	87	0.00
59 T	2-Hexanone	0.372	0.353	5.1	86	0.00
60 T	Dibromochloromethane	0.419	0.398	5.0	91	0.00
61 T	1,2-Dibromoethane	0.367	0.338	7.9	88	0.00
62 S	4-Bromofluorobenzene	0.416	0.415	0.2	97	0.00
63 I	Chlorobenzene-d5	1.000	1.000	0.0	91	0.00
64 T	Tetrachloroethene	0.364	0.332	8.8	86	0.00
65 PM	Chlorobenzene	1.094	1.057	3.4	88	0.00
66 T	1,1,1,2-Tetrachloroethane	0.391	0.388	0.8	93	0.00
67 C	Ethyl Benzene	1.786	1.788	-0.1#	86	0.00
68 T	m/p-Xylenes	0.696	0.717	-3.0	91	0.00
69 T	o-Xylene	0.652	0.672	-3.1	90	0.00
70 T	Styrene	1.119	1.151	-2.9	90	0.00
71 P	Bromoform	0.349	0.359	-2.9	100	0.00
72 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	97	0.00
73 T	Isopropylbenzene	3.369	3.604	-7.0	96	0.00
74 T	N-ethyl acetate	1.464	1.724	-17.8	96	0.00
75 P	1,1,2,2-Tetrachloroethane	1.137	1.212	-6.6	100	0.00
76 T	1,2,3-Trichloropropane	0.931	0.996	-7.0	101	0.00
77 T	Bromobenzene	0.899	0.906	-0.8	99	0.00
78 T	n-propylbenzene	3.902	4.092	-4.9	92	0.00
79 T	2-Chlorotoluene	2.380	2.451	-3.0	94	0.00
80 T	1,3,5-Trimethylbenzene	2.770	2.875	-3.8	95	0.00
81 T	trans-1,4-Dichloro-2-butene	0.392	0.422	-7.7	99	0.00
82 T	4-Chlorotoluene	2.777	2.916	-5.0	95	0.00
83 T	tert-Butylbenzene	2.878	2.886	-0.3	93	0.00
84 T	1,2,4-Trimethylbenzene	2.794	2.883	-3.2	94	0.00
85 T	sec-Butylbenzene	3.488	3.530	-1.2	91	0.00
86 T	p-Isopropyltoluene	2.981	3.001	-0.7	92	0.00
87 T	1,3-Dichlorobenzene	1.658	1.626	1.9	96	0.00
88 T	1,4-Dichlorobenzene	1.713	1.637	4.4	96	0.00
89 T	n-Butylbenzene	2.791	2.833	-1.5	93	0.00
90 T	Hexachloroethane	0.608	0.571	6.1	93	0.00
91 T	1,2-Dichlorobenzene	1.607	1.546	3.8	95	0.00
92 T	1,2-Dibromo-3-Chloropropane	0.277	0.281	-1.4	102	0.00
93 T	1,2,4-Trichlorobenzene	1.062	0.998	6.0	97	0.00
94 T	Hexachlorobutadiene	0.439	0.396	9.8	97	0.00
95 T	Naphthalene	3.187	3.414	-7.1	105	0.00

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120425\
Data File : VX048721.D
Acq On : 04 Dec 2025 14:04
Operator : JC/MD
Sample : VSTDICV050
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
ICVVX120425

Quant Time: Dec 05 03:31:44 2025
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X120425W.M
Quant Title : SW846 8260
QLast Update : Fri Dec 05 03:27:34 2025
Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
Max. RRF Dev : 25% Max. Rel. Area : 150%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
96 T	1,2,3-Trichlorobenzene	0.979	0.967	1.2	101	0.00

(#) = Out of Range

SPCC's out = 0 CCC's out = 6

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\WX120425\
 Data File : VX048721.D
 Acq On : 04 Dec 2025 14:04
 Operator : JC/MD
 Sample : VSTDICV050
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 ICVVX120425

Quant Time: Dec 05 03:31:44 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X120425W.M
 Quant Title : SW846 8260
 QLast Update : Fri Dec 05 03:27:34 2025
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
 Max. RRF Dev : 25% Max. Rel. Area : 150%

	Compound	Amount	Calc.	%Dev	Area%	Dev(min)
1 I	Pentafluorobenzene	50.000	50.000	0.0	99	0.00
2 T	Dichlorodifluoromethane	50.000	58.738	-17.5	104	0.00
3 P	Chloromethane	50.000	48.826	2.3	97	0.00
4 C	Vinyl Chloride	50.000	48.559	2.9#	93	0.00
5 T	Bromomethane	50.000	54.893	-9.8	99	0.00
6 T	Chloroethane	50.000	48.675	2.7	97	0.00
7 T	Trichlorofluoromethane	50.000	48.801	2.4	92	0.00
8 T	Diethyl Ether	50.000	48.151	3.7	98	0.00
9 T	1,1,2-Trichlorotrifluoroeth	50.000	45.717	8.6	87	0.00
10 T	Methyl Iodide	50.000	50.395	-0.8	94	0.00
11 T	Tert butyl alcohol	250.000	250.243	-0.1	94	0.00
12 CM	1,1-Dichloroethene	50.000	46.649	6.7#	89	0.00
13 T	Acrolein	250.000	232.010	7.2	111	0.00
14 T	Allyl chloride	50.000	42.572	14.9	77	0.00
15 T	Acrylonitrile	250.000	226.688	9.3	83	0.00
16 T	Acetone	250.000	204.537	18.2	85	0.00
17 T	Carbon Disulfide	50.000	41.077	17.8	77	0.00
18 T	Methyl Acetate	50.000	44.401	11.2	82	0.00
19 T	Methyl tert-butyl Ether	50.000	45.629	8.7	86	0.00
20 T	Methylene Chloride	50.000	47.622	4.8	93	0.00
21 T	trans-1,2-Dichloroethene	50.000	44.308	11.4	83	0.00
22 T	Diisopropyl ether	50.000	50.527	-1.1	116	0.00
23 T	Vinyl Acetate	250.000	276.044	-10.4	121	0.00
24 P	1,1-Dichloroethane	50.000	52.811	-5.6	117	0.00
25 T	2-Butanone	250.000	244.061	2.4	109	0.00
26 T	2,2-Dichloropropane	50.000	45.857	8.3	98	0.00
27 T	cis-1,2-Dichloroethene	50.000	48.033	3.9	101	0.00
28 T	Bromochloromethane	50.000	45.040	9.9	107	0.00
29 T	Tetrahydrofuran	250.000	277.876	-11.2	129	0.00
30 C	Chloroform	50.000	52.061	-4.1#	115	0.00
31 T	Cyclohexane	50.000	49.171	1.7	111	0.00
32 T	1,1,1-Trichloroethane	50.000	50.071	-0.1	106	0.00
33 S	1,2-Dichloroethane-d4	50.000	54.154	-8.3	125	0.00
34 I	1,4-Difluorobenzene	50.000	50.000	0.0	107	0.00
35 S	Dibromofluoromethane	50.000	49.960	0.1	110	0.00
36 T	1,1-Dichloropropene	50.000	47.458	5.1	110	0.00
37 T	Ethyl Acetate	50.000	46.283	7.4	104	0.00
38 T	Carbon Tetrachloride	50.000	47.169	5.7	102	0.00
39 T	Methylcyclohexane	50.000	42.440	15.1	79	0.00
40 TM	Benzene	50.000	50.374	-0.7	114	0.00
41 T	Methacrylonitrile	50.000	48.603	2.8	110	0.00
42 TM	1,2-Dichloroethane	50.000	51.045	-2.1	120	0.00
43 T	Isopropyl Acetate	50.000	53.729	-7.5	130	0.00
44 TM	Trichloroethene	50.000	45.250	9.5	97	0.00
45 C	1,2-Dichloropropane	50.000	42.935	14.1#	84	0.00
46 T	Dibromomethane	50.000	43.757	12.5	86	0.00
47 T	Bromodichloromethane	50.000	42.902	14.2	87	0.00
48 T	Methyl methacrylate	50.000	42.366	15.3	84	0.00

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120425\
 Data File : VX048721.D
 Acq On : 04 Dec 2025 14:04
 Operator : JC/MD
 Sample : VSTDICV050
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 ICVVX120425

Quant Time: Dec 05 03:31:44 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X120425W.M
 Quant Title : SW846 8260
 QLast Update : Fri Dec 05 03:27:34 2025
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
 Max. RRF Dev : 25% Max. Rel. Area : 150%

	Compound	Amount	Calc.	%Dev	Area%	Dev(min)
49 T	1,4-Dioxane	1000.000	985.778	1.4	99	0.00
50 S	Toluene-d8	50.000	44.544	10.9	88	0.00
51 T	4-Methyl-2-Pentanone	250.000	220.496	11.8	83	0.00
52 CM	Toluene	50.000	46.077	7.8#	86	0.00
53 T	t-1,3-Dichloropropene	50.000	46.514	7.0	87	0.00
54 T	cis-1,3-Dichloropropene	50.000	45.699	8.6	86	0.00
55 T	1,1,2-Trichloroethane	50.000	45.811	8.4	88	0.00
56 T	Ethyl methacrylate	50.000	48.247	3.5	88	0.00
57 T	1,3-Dichloropropane	50.000	45.415	9.2	88	0.00
58 T	2-Chloroethyl Vinyl ether	250.000	242.598	3.0	87	0.00
59 T	2-Hexanone	250.000	237.637	4.9	86	0.00
60 T	Dibromochloromethane	50.000	47.409	5.2	91	0.00
61 T	1,2-Dibromoethane	50.000	46.076	7.8	88	0.00
62 S	4-Bromofluorobenzene	50.000	49.905	0.2	97	0.00
63 I	Chlorobenzene-d5	50.000	50.000	0.0	91	0.00
64 T	Tetrachloroethene	50.000	45.571	8.9	86	0.00
65 PM	Chlorobenzene	50.000	48.287	3.4	88	0.00
66 T	1,1,1,2-Tetrachloroethane	50.000	49.615	0.8	93	0.00
67 C	Ethyl Benzene	50.000	50.069	-0.1#	86	0.00
68 T	m/p-Xylenes	100.000	103.051	-3.1	91	0.00
69 T	o-Xylene	50.000	51.580	-3.2	90	0.00
70 T	Styrene	50.000	51.437	-2.9	90	0.00
71 P	Bromoform	50.000	51.388	-2.8	100	0.00
72 I	1,4-Dichlorobenzene-d4	50.000	50.000	0.0	97	0.00
73 T	Isopropylbenzene	50.000	53.483	-7.0	96	0.00
74 T	N-amyl acetate	50.000	58.895	-17.8	96	0.00
75 P	1,1,2,2-Tetrachloroethane	50.000	53.282	-6.6	100	0.00
76 T	1,2,3-Trichloropropane	50.000	53.497	-7.0	101	0.00
77 T	Bromobenzene	50.000	50.357	-0.7	99	0.00
78 T	n-propylbenzene	50.000	52.425	-4.8	92	0.00
79 T	2-Chlorotoluene	50.000	51.501	-3.0	94	0.00
80 T	1,3,5-Trimethylbenzene	50.000	51.894	-3.8	95	0.00
81 T	trans-1,4-Dichloro-2-butene	50.000	53.840	-7.7	99	0.00
82 T	4-Chlorotoluene	50.000	52.505	-5.0	95	0.00
83 T	tert-Butylbenzene	50.000	50.138	-0.3	93	0.00
84 T	1,2,4-Trimethylbenzene	50.000	51.592	-3.2	94	0.00
85 T	sec-Butylbenzene	50.000	50.606	-1.2	91	0.00
86 T	p-Isopropyltoluene	50.000	50.326	-0.7	92	0.00
87 T	1,3-Dichlorobenzene	50.000	49.048	1.9	96	0.00
88 T	1,4-Dichlorobenzene	50.000	47.783	4.4	96	0.00
89 T	n-Butylbenzene	50.000	50.745	-1.5	93	0.00
90 T	Hexachloroethane	50.000	46.926	6.1	93	0.00
91 T	1,2-Dichlorobenzene	50.000	48.097	3.8	95	0.00
92 T	1,2-Dibromo-3-Chloropropane	50.000	50.715	-1.4	102	0.00
93 T	1,2,4-Trichlorobenzene	50.000	46.996	6.0	97	0.00
94 T	Hexachlorobutadiene	50.000	45.101	9.8	97	0.00
95 T	Naphthalene	50.000	55.628	-11.3	105	0.00

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120425\
Data File : VX048721.D
Acq On : 04 Dec 2025 14:04
Operator : JC/MD
Sample : VSTDICV050
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
ICVVX120425

Quant Time: Dec 05 03:31:44 2025
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X120425W.M
Quant Title : SW846 8260
QLast Update : Fri Dec 05 03:27:34 2025
Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
Max. RRF Dev : 25% Max. Rel. Area : 150%

	Compound	Amount	Calc.	%Dev	Area%	Dev(min)
96 T	1,2,3-Trichlorobenzene	50.000	49.386	1.2	101	0.00

(#) = Out of Range SPCC's out = 0 CCC's out = 6



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

VOLATILE CONTINUING CALIBRATION CHECK

Lab Name:	<u>Alliance</u>	Contract:	<u>JAC005</u>
Lab Code:	<u>ACE</u>	SDG No.:	<u>Q3757</u>
Instrument ID:	<u>MSVOA_N</u>	Calibration Date/Time:	<u>12/04/2025</u> <u>11:44</u>
Lab File ID:	<u>VN088437.D</u>	Init. Calib. Date(s):	<u>11/24/2025</u> <u>11/24/2025</u>
Heated Purge: (Y/N)	<u>N</u>	Init. Calib. Time(s):	<u>12:38</u> <u>14:35</u>
GC Column:	<u>RXI-624</u>	ID:	<u>0.25</u> (mm)

COMPOUND	RRF	RRF050	MIN RRF	%D	MAX%D
Vinyl Chloride	0.821	0.817		-0.49	20
1,1-Dichloroethene	0.610	0.619		1.48	20
1,1-Dichloroethane	1.367	1.487	0.1	8.78	20
cis-1,2-Dichloroethene	0.785	0.832		5.99	20
1,1,1-Trichloroethane	1.181	1.277		8.13	20
Benzene	1.570	1.743		11.02	20
1,2-Dichloroethane	0.599	0.729		21.7	20
Trichloroethene	0.371	0.385		3.77	20
1,1,2-Trichloroethane	0.351	0.402		14.53	20
Tetrachloroethene	0.396	0.360		-9.09	20
1,2-Dichloroethane-d4	0.942	1.041		10.51	20
Dibromofluoromethane	0.346	0.396		14.45	20
Toluene-d8	1.289	1.429		10.86	20
4-Bromofluorobenzene	0.442	0.507		14.71	20

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_N\Data\VN120425\
 Data File : VN088437.D
 Acq On : 04 Dec 2025 11:44
 Operator : JC\MD
 Sample : VSTDCCC050
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 VSTDCCC050

Manual Integrations
 APPROVED

Reviewed By : John Carlone 12/05/2025
 Supervised By : Semsettin Yesilyurt 12/05/2025

Quant Time: Dec 05 00:11:19 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N112425W.M
 Quant Title : SW846 8260
 QLast Update : Tue Nov 25 01:34:52 2025
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.206	168	259230	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.077	114	454552	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.841	117	431157	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.765	152	210277	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.559	65	269836	55.256	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery = 110.520%			
35) Dibromofluoromethane	8.147	113	180201	57.206	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery = 114.420%			
50) Toluene-d8	10.541	98	649537	55.419	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery = 110.840%			
62) 4-Bromofluorobenzene	12.829	95	230283	57.262	ug/l	0.00
Spiked Amount 50.000	Range 77 - 121		Recovery = 114.520%			
Target Compounds					Qvalue	
2) Dichlorodifluoromethane	2.142	85	201478	51.745	ug/l	99
3) Chloromethane	2.383	50	206193	49.307	ug/l	96
4) Vinyl Chloride	2.536	62	211796	49.784	ug/l	100
5) Bromomethane	2.977	94	100901	49.278	ug/l	93
6) Chloroethane	3.142	64	138625	52.654	ug/l	95
7) Trichlorofluoromethane	3.512	101	324899	54.337	ug/l	99
8) Diethyl Ether	3.959	74	117687	52.439	ug/l	99
9) 1,1,2-Trichlorotrifluo...	4.365	101	164403	51.647	ug/l	99
10) Methyl Iodide	4.583	142	141417	44.878	ug/l	92
11) Tert butyl alcohol	5.524	59	174143	223.357	ug/l	98
12) 1,1-Dichloroethene	4.336	96	160467	50.738	ug/l	93
13) Acrolein	4.171	56	216629	290.220	ug/l	98
14) Allyl chloride	5.018	41	291153	49.668	ug/l	96
15) Acrylonitrile	5.706	53	590127	262.341	ug/l	99
16) Acetone	4.418	43	408907	231.533	ug/l	100
17) Carbon Disulfide	4.706	76	513890	49.724	ug/l	99
18) Methyl Acetate	5.012	43	289409	55.051	ug/l	99
19) Methyl tert-butyl Ether	5.789	73	650434	54.892	ug/l	95
20) Methylene Chloride	5.265	84	193652	51.250	ug/l	95
21) trans-1,2-Dichloroethene	5.771	96	182717	52.153	ug/l	99
22) Diisopropyl ether	6.659	45	700009	55.934	ug/l	98
23) Vinyl Acetate	6.589	43	3136556	311.072	ug/l	99
24) 1,1-Dichloroethane	6.553	63	385388	54.378	ug/l	99
25) 2-Butanone	7.465	43	739812	260.572	ug/l	98
26) 2,2-Dichloropropane	7.471	77	326583	53.386	ug/l	99
27) cis-1,2-Dichloroethene	7.465	96	215724	52.983	ug/l	99
28) Bromochloromethane	7.794	49	199532	56.516	ug/l	97
29) Tetrahydrofuran	7.818	42	562734	276.329	ug/l	99
30) Chloroform	7.947	83	392592	56.375	ug/l	99
31) Cyclohexane	8.241	56	322585	50.228	ug/l	99
32) 1,1,1-Trichloroethane	8.147	97	331052	54.066	ug/l	99
36) 1,1-Dichloropropene	8.347	75	262020	53.897	ug/l	99
37) Ethyl Acetate	7.542	43	348681	54.547	ug/l	100
38) Carbon Tetrachloride	8.341	117	280878	58.038	ug/l	96
39) Methylcyclohexane	9.577	83	272178	52.341	ug/l	94
40) Benzene	8.583	78	792340	55.531	ug/l	98

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN120425\
 Data File : VN088437.D
 Acq On : 04 Dec 2025 11:44
 Operator : JC\MD
 Sample : VSTDCCC050
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 VSTDCCC050

Manual Integrations
 APPROVED

Reviewed By :John Carlone 12/05/2025
 Supervised By :Semsettin Yesilyurt 12/05/2025

Quant Time: Dec 05 00:11:19 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N112425W.M
 Quant Title : SW846 8260
 QLast Update : Tue Nov 25 01:34:52 2025
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Methacrylonitrile	7.759	41	184821	58.376	ug/l	96
42) 1,2-Dichloroethane	8.647	62	331550	60.904	ug/l	97
43) Isopropyl Acetate	8.671	43	542837	55.581	ug/l	100
44) Trichloroethene	9.330	130	174795	51.846	ug/l	96
45) 1,2-Dichloropropane	9.594	63	210122	57.477	ug/l	96
46) Dibromomethane	9.688	93	150122	57.541	ug/l	98
47) Bromodichloromethane	9.865	83	307416	57.749	ug/l	97
48) Methyl methacrylate	9.659	41	262024	58.800	ug/l	97
49) 1,4-Dioxane	9.677	88	58650	1034.918	ug/l #	97
51) 4-Methyl-2-Pentanone	10.424	43	1723595	299.404	ug/l	100
52) Toluene	10.606	92	480292	58.295	ug/l	98
53) t-1,3-Dichloropropene	10.812	75	312027	57.105	ug/l	98
54) cis-1,3-Dichloropropene	10.288	75	330756	57.561	ug/l	97
55) 1,1,2-Trichloroethane	10.994	97	182938	57.390	ug/l	98
56) Ethyl methacrylate	10.853	69	311735	61.911	ug/l	98
57) 1,3-Dichloropropane	11.141	76	332821	57.876	ug/l	99
58) 2-Chloroethyl Vinyl ether	10.135	63	883178	304.242	ug/l	100
59) 2-Hexanone	11.177	43	1099155	303.090	ug/l	100
60) Dibromochloromethane	11.335	129	214567	59.046	ug/l	99
61) 1,2-Dibromoethane	11.447	107	184522	57.164	ug/l	99
64) Tetrachloroethene	11.077	164	155230	45.415	ug/l	97
65) Chlorobenzene	11.871	112	504973	52.142	ug/l	99
66) 1,1,1,2-Tetrachloroethane	11.935	131	177636	54.384	ug/l	98
67) Ethyl Benzene	11.941	91	910559	53.794	ug/l	100
68) m/p-Xylenes	12.047	106	690769	110.067	ug/l	100
69) o-Xylene	12.376	106	319328	53.371	ug/l	97
70) Styrene	12.388	104	558388	57.477	ug/l	99
71) Bromoform	12.559	173	134945	56.405	ug/l #	99
73) Isopropylbenzene	12.671	105	846156	54.424	ug/l	100
74) N-amyl acetate	12.500	43	212974m	40.223	ug/l	
75) 1,1,2,2-Tetrachloroethane	12.918	83	276796	55.070	ug/l	98
76) 1,2,3-Trichloropropane	12.971	75	264559m	53.717	ug/l	
77) Bromobenzene	12.959	156	197041	55.390	ug/l	98
78) n-propylbenzene	13.012	91	1046093	54.998	ug/l	99
79) 2-Chlorotoluene	13.100	91	615186	52.964	ug/l	99
80) 1,3,5-Trimethylbenzene	13.147	105	720007	55.887	ug/l	98
81) trans-1,4-Dichloro-2-b...	12.712	75	100456	57.865	ug/l #	89
82) 4-Chlorotoluene	13.200	91	656816	56.933	ug/l	98
83) tert-Butylbenzene	13.412	119	588225	53.710	ug/l	99
84) 1,2,4-Trimethylbenzene	13.459	105	730000	57.007	ug/l	98
85) sec-Butylbenzene	13.588	105	847596	53.872	ug/l	99
86) p-Isopropyltoluene	13.706	119	706011	54.723	ug/l	100
87) 1,3-Dichlorobenzene	13.712	146	376351	54.348	ug/l	99
88) 1,4-Dichlorobenzene	13.788	146	380675	52.479	ug/l	97
89) n-Butylbenzene	14.029	91	671306	53.590	ug/l	100
90) Hexachloroethane	14.306	117	127576	54.130	ug/l	98
91) 1,2-Dichlorobenzene	14.082	146	349875	53.690	ug/l	99
92) 1,2-Dibromo-3-Chloropr...	14.694	75	63936	52.428	ug/l	96
93) 1,2,4-Trichlorobenzene	15.365	180	194146	50.142	ug/l	98
94) Hexachlorobutadiene	15.470	225	72189	51.883	ug/l	97
95) Naphthalene	15.612	128	712709	53.119	ug/l	99
96) 1,2,3-Trichlorobenzene	15.812	180	192507	51.087	ug/l	97

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN120425\
Data File : VN088437.D
Acq On : 04 Dec 2025 11:44
Operator : JC\MD
Sample : VSTDCCC050
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 3 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
VSTDCCC050

Manual Integrations
APPROVED

Reviewed By :John Carlone 12/05/2025
Supervised By :Semsettin Yesilyurt 12/05/2025

Quant Time: Dec 05 00:11:19 2025
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N112425W.M
Quant Title : SW846 8260
QLast Update : Tue Nov 25 01:34:52 2025
Response via : Initial Calibration

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

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN120425\
Data File : VN088437.D
Acq On : 04 Dec 2025 11:44
Operator : JC\MD
Sample : VSTDCCC050
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 3 Sample Multiplier: 1

Instrument :

MSVOA_N

ClientSampleId :

VSTDCCC050

Manual Integrations

APPROVED

Reviewed By :John Carlone 12/05/2025

Supervised By :Semsettin Yesilyurt 12/05/2025

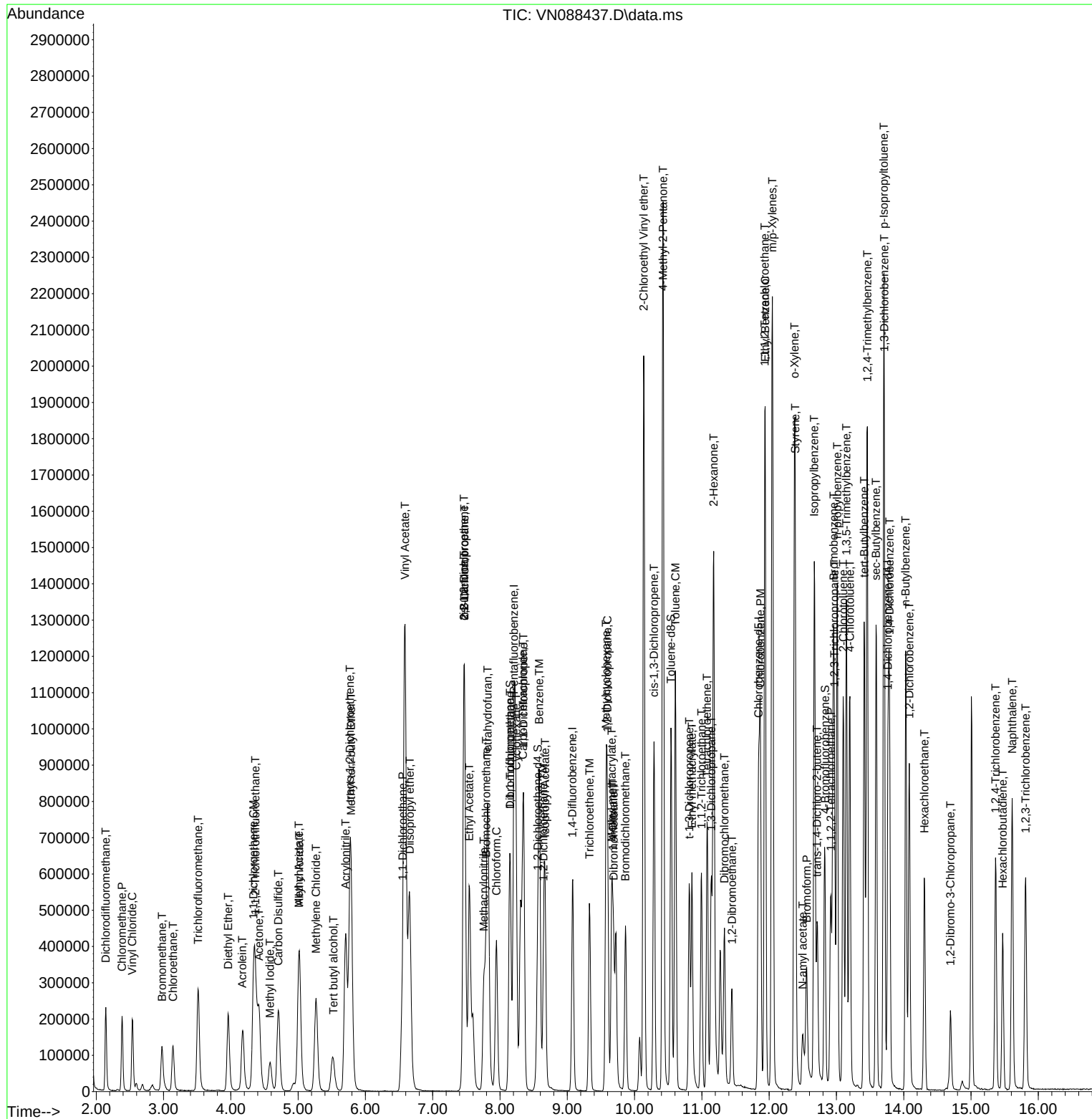
Quant Time: Dec 05 00:11:19 2025

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N112425W.M

Quant Title : SW846 8260

QLast Update : Tue Nov 25 01:34:52 2025

Response via : Initial Calibration



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN120425\
 Data File : VN088437.D
 Acq On : 04 Dec 2025 11:44
 Operator : JC\MD
 Sample : VSTDCCC050
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 MSVOA_N
 LabSampleId :
 VSTDCCC050

Quant Time: Dec 05 00:11:19 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N112425W.M
 Quant Title : SW846 8260
 QLast Update : Tue Nov 25 01:34:52 2025
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
 Max. RRF Dev : 25% Max. Rel. Area : 150%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
1 I	Pentafluorobenzene	1.000	1.000	0.0	102	0.00
2 T	Dichlorodifluoromethane	0.751	0.777	-3.5	107	0.00
3 P	Chloromethane	0.807	0.795	1.5	101	0.00
4 C	Vinyl Chloride	0.821	0.817	0.5#	103	0.00
5 T	Bromomethane	0.395	0.389	1.5	100	0.00
6 T	Chloroethane	0.508	0.535	-5.3	108	0.00
7 T	Trichlorofluoromethane	1.153	1.253	-8.7	117	0.00
8 T	Diethyl Ether	0.433	0.454	-4.8	104	0.00
9 T	1,1,2-Trichlorotrifluoroeth	0.614	0.634	-3.3	115	0.00
10 T	Methyl Iodide	0.608	0.546	10.2	84	0.00
11 T	Tert butyl alcohol	0.150	0.134	10.7	85	0.01
12 CM	1,1-Dichloroethene	0.610	0.619	-1.5#	111	0.00
13 T	Acrolein	0.144	0.167	-16.0	113	0.00
14 T	Allyl chloride	1.131	1.123	0.7	100	0.00
15 T	Acrylonitrile	0.434	0.455	-4.8	99	0.00
16 T	Acetone	0.341	0.315	7.6	96	0.00
17 T	Carbon Disulfide	1.993	1.982	0.6	102	0.00
18 T	Methyl Acetate	1.014	1.116	-10.1	108	0.00
19 T	Methyl tert-butyl Ether	2.285	2.509	-9.8	102	0.00
20 T	Methylene Chloride	0.729	0.747	-2.5	101	0.00
21 T	trans-1,2-Dichloroethene	0.676	0.705	-4.3	105	0.00
22 T	Diisopropyl ether	2.414	2.700	-11.8	107	0.00
23 T	Vinyl Acetate	1.945	2.420	-24.4	115	0.00
24 P	1,1-Dichloroethane	1.367	1.487	-8.8	106	0.00
25 T	2-Butanone	0.548	0.571	-4.2	98	0.00
26 T	2,2-Dichloropropane	1.180	1.260	-6.8	107	0.00
27 T	cis-1,2-Dichloroethene	0.785	0.832	-6.0	102	0.00
28 T	Bromochloromethane	0.681	0.770	-13.1	107	0.00
29 T	Tetrahydrofuran	0.393	0.434	-10.4	103	0.00
30 C	Chloroform	1.343	1.514	-12.7#	111	0.00
31 T	Cyclohexane	1.239	1.244	-0.4	112	0.00
32 T	1,1,1-Trichloroethane	1.181	1.277	-8.1	112	0.00
33 S	1,2-Dichloroethane-d4	0.942	1.041	-10.5	106	0.00
34 I	1,4-Difluorobenzene	1.000	1.000	0.0	95	0.00
35 S	Dibromofluoromethane	0.346	0.396	-14.5	107	0.00
36 T	1,1-Dichloropropene	0.535	0.576	-7.7	111	0.00
37 T	Ethyl Acetate	0.703	0.767	-9.1	101	0.00
38 T	Carbon Tetrachloride	0.532	0.618	-16.2	117	0.00
39 T	Methylcyclohexane	0.572	0.599	-4.7	107	0.00
40 TM	Benzene	1.570	1.743	-11.0	106	0.00
41 T	Methacrylonitrile	0.348	0.407	-17.0	108	0.00
42 TM	1,2-Dichloroethane	0.599	0.729	-21.7	112	0.00
43 T	Isopropyl Acetate	1.074	1.194	-11.2	101	0.00
44 TM	Trichloroethene	0.371	0.385	-3.8	102	0.00
45 C	1,2-Dichloropropane	0.402	0.462	-14.9#	108	0.00
46 T	Dibromomethane	0.287	0.330	-15.0	107	0.00
47 T	Bromodichloromethane	0.586	0.676	-15.4	108	0.00
48 T	Methyl methacrylate	0.490	0.576	-17.6	106	0.00

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN120425\
 Data File : VN088437.D
 Acq On : 04 Dec 2025 11:44
 Operator : JC\MD
 Sample : VSTDCCC050
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 MSVOA_N
 LabSampleId :
 VSTDCCC050

Quant Time: Dec 05 00:11:19 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N112425W.M
 Quant Title : SW846 8260
 QLast Update : Tue Nov 25 01:34:52 2025
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
 Max. RRF Dev : 25% Max. Rel. Area : 150%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
49 T	1,4-Dioxane	0.006	0.006	0.0	92	0.00
50 S	Toluene-d8	1.289	1.429	-10.9	105	0.00
51 T	4-Methyl-2-Pentanone	0.633	0.758	-19.7	107	0.00
52 CM	Toluene	0.906	1.057	-16.7#	109	0.00
53 T	t-1,3-Dichloropropene	0.601	0.686	-14.1	103	0.00
54 T	cis-1,3-Dichloropropene	0.632	0.728	-15.2	104	0.00
55 T	1,1,2-Trichloroethane	0.351	0.402	-14.5	108	0.00
56 T	Ethyl methacrylate	0.554	0.686	-23.8	105	0.00
57 T	1,3-Dichloropropane	0.633	0.732	-15.6	107	0.00
58 T	2-Chloroethyl Vinyl ether	0.319	0.389	-21.9	101	0.00
59 T	2-Hexanone	0.399	0.484	-21.3	105	0.00
60 T	Dibromochloromethane	0.400	0.472	-18.0	109	0.00
61 T	1,2-Dibromoethane	0.355	0.406	-14.4	104	0.00
62 S	4-Bromofluorobenzene	0.442	0.507	-14.7	107	0.00
63 I	Chlorobenzene-d5	1.000	1.000	0.0	101	0.00
64 T	Tetrachloroethene	0.396	0.360	9.1	97	0.00
65 PM	Chlorobenzene	1.123	1.171	-4.3	107	0.00
66 T	1,1,1,2-Tetrachloroethane	0.379	0.412	-8.7	109	0.00
67 C	Ethyl Benzene	1.963	2.112	-7.6#	110	0.00
68 T	m/p-Xylenes	0.728	0.801	-10.0	111	0.00
69 T	o-Xylene	0.694	0.741	-6.8	108	0.00
70 T	Styrene	1.127	1.295	-14.9	110	0.00
71 P	Bromoform	0.277	0.313	-13.0	111	0.00
72 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	101	0.00
73 T	Isopropylbenzene	3.697	4.024	-8.8	112	0.00
74 T	N-amyl acetate	1.259	1.013	19.5	98	0.00
75 P	1,1,2,2-Tetrachloroethane	1.195	1.316	-10.1	118	0.00
76 T	1,2,3-Trichloropropane	1.171	1.258	-7.4	106	0.00
77 T	Bromobenzene	0.846	0.937	-10.8	109	0.00
78 T	n-propylbenzene	4.523	4.975	-10.0	113	0.00
79 T	2-Chlorotoluene	2.762	2.926	-5.9	109	0.00
80 T	1,3,5-Trimethylbenzene	3.063	3.424	-11.8	113	0.00
81 T	trans-1,4-Dichloro-2-butene	0.413	0.478	-15.7	112	0.00
82 T	4-Chlorotoluene	2.743	3.124	-13.9	110	0.00
83 T	tert-Butylbenzene	2.604	2.797	-7.4	112	0.00
84 T	1,2,4-Trimethylbenzene	3.045	3.472	-14.0	112	0.00
85 T	sec-Butylbenzene	3.741	4.031	-7.8	113	0.00
86 T	p-Isopropyltoluene	3.068	3.358	-9.5	111	0.00
87 T	1,3-Dichlorobenzene	1.647	1.790	-8.7	112	0.00
88 T	1,4-Dichlorobenzene	1.725	1.810	-4.9	109	0.00
89 T	n-Butylbenzene	2.979	3.192	-7.2	112	0.00
90 T	Hexachloroethane	0.560	0.607	-8.4	114	0.00
91 T	1,2-Dichlorobenzene	1.550	1.664	-7.4	110	0.00
92 T	1,2-Dibromo-3-Chloropropane	0.290	0.304	-4.8	106	0.00
93 T	1,2,4-Trichlorobenzene	0.921	0.923	-0.2	105	0.00
94 T	Hexachlorobutadiene	0.331	0.343	-3.6	117	0.00
95 T	Naphthalene	3.190	3.389	-6.2	104	0.00

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN120425\
Data File : VN088437.D
Acq On : 04 Dec 2025 11:44
Operator : JC\MD
Sample : VSTDCCC050
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 3 Sample Multiplier: 1

Instrument :
MSVOA_N
LabSampleId :
VSTDCCC050

Quant Time: Dec 05 00:11:19 2025
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N112425W.M
Quant Title : SW846 8260
QLast Update : Tue Nov 25 01:34:52 2025
Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
Max. RRF Dev : 25% Max. Rel. Area : 150%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
96 T	1,2,3-Trichlorobenzene	0.896	0.915	-2.1	105	0.00

(#) = Out of Range

SPCC's out = 0 CCC's out = 6

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN120425\
 Data File : VN088437.D
 Acq On : 04 Dec 2025 11:44
 Operator : JC\MD
 Sample : VSTDCCC050
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 MSVOA_N
 LabSampleId :
 VSTDCCC050

Quant Time: Dec 05 00:11:19 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N112425W.M
 Quant Title : SW846 8260
 QLast Update : Tue Nov 25 01:34:52 2025
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
 Max. RRF Dev : 25% Max. Rel. Area : 150%

	Compound	Amount	Calc.	%Dev	Area%	Dev(min)
1 I	Pentafluorobenzene	50.000	50.000	0.0	102	0.00
2 T	Dichlorodifluoromethane	50.000	51.745	-3.5	107	0.00
3 P	Chloromethane	50.000	49.307	1.4	101	0.00
4 C	Vinyl Chloride	50.000	49.784	0.4#	103	0.00
5 T	Bromomethane	50.000	49.278	1.4	100	0.00
6 T	Chloroethane	50.000	52.654	-5.3	108	0.00
7 T	Trichlorofluoromethane	50.000	54.337	-8.7	117	0.00
8 T	Diethyl Ether	50.000	52.439	-4.9	104	0.00
9 T	1,1,2-Trichlorotrifluoroeth	50.000	51.647	-3.3	115	0.00
10 T	Methyl Iodide	50.000	44.878	10.2	84	0.00
11 T	Tert butyl alcohol	250.000	223.357	10.7	85	0.01
12 CM	1,1-Dichloroethene	50.000	50.738	-1.5#	111	0.00
13 T	Acrolein	250.000	290.220	-16.1	113	0.00
14 T	Allyl chloride	50.000	49.668	0.7	100	0.00
15 T	Acrylonitrile	250.000	262.341	-4.9	99	0.00
16 T	Acetone	250.000	231.533	7.4	96	0.00
17 T	Carbon Disulfide	50.000	49.724	0.6	102	0.00
18 T	Methyl Acetate	50.000	55.051	-10.1	108	0.00
19 T	Methyl tert-butyl Ether	50.000	54.892	-9.8	102	0.00
20 T	Methylene Chloride	50.000	51.250	-2.5	101	0.00
21 T	trans-1,2-Dichloroethene	50.000	52.153	-4.3	105	0.00
22 T	Diisopropyl ether	50.000	55.934	-11.9	107	0.00
23 T	Vinyl Acetate	250.000	311.072	-24.4	115	0.00
24 P	1,1-Dichloroethane	50.000	54.378	-8.8	106	0.00
25 T	2-Butanone	250.000	260.572	-4.2	98	0.00
26 T	2,2-Dichloropropane	50.000	53.386	-6.8	107	0.00
27 T	cis-1,2-Dichloroethene	50.000	52.983	-6.0	102	0.00
28 T	Bromochloromethane	50.000	56.516	-13.0	107	0.00
29 T	Tetrahydrofuran	250.000	276.329	-10.5	103	0.00
30 C	Chloroform	50.000	56.375	-12.8#	111	0.00
31 T	Cyclohexane	50.000	50.228	-0.5	112	0.00
32 T	1,1,1-Trichloroethane	50.000	54.066	-8.1	112	0.00
33 S	1,2-Dichloroethane-d4	50.000	55.256	-10.5	106	0.00
34 I	1,4-Difluorobenzene	50.000	50.000	0.0	95	0.00
35 S	Dibromofluoromethane	50.000	57.206	-14.4	107	0.00
36 T	1,1-Dichloropropene	50.000	53.897	-7.8	111	0.00
37 T	Ethyl Acetate	50.000	54.547	-9.1	101	0.00
38 T	Carbon Tetrachloride	50.000	58.038	-16.1	117	0.00
39 T	Methylcyclohexane	50.000	52.341	-4.7	107	0.00
40 TM	Benzene	50.000	55.531	-11.1	106	0.00
41 T	Methacrylonitrile	50.000	58.376	-16.8	108	0.00
42 TM	1,2-Dichloroethane	50.000	60.904	-21.8	112	0.00
43 T	Isopropyl Acetate	50.000	55.581	-11.2	101	0.00
44 TM	Trichloroethene	50.000	51.846	-3.7	102	0.00
45 C	1,2-Dichloropropane	50.000	57.477	-15.0#	108	0.00
46 T	Dibromomethane	50.000	57.541	-15.1	107	0.00
47 T	Bromodichloromethane	50.000	57.749	-15.5	108	0.00
48 T	Methyl methacrylate	50.000	58.800	-17.6	106	0.00

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN120425\
 Data File : VN088437.D
 Acq On : 04 Dec 2025 11:44
 Operator : JC\MD
 Sample : VSTDCCC050
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 MSVOA_N
 LabSampleId :
 VSTDCCC050

Quant Time: Dec 05 00:11:19 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N112425W.M
 Quant Title : SW846 8260
 QLast Update : Tue Nov 25 01:34:52 2025
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
 Max. RRF Dev : 25% Max. Rel. Area : 150%

	Compound	Amount	Calc.	%Dev	Area%	Dev(min)

49 T	1,4-Dioxane	1000.000	1034.918	-3.5	92	0.00
50 S	Toluene-d8	50.000	55.419	-10.8	105	0.00
51 T	4-Methyl-2-Pentanone	250.000	299.404	-19.8	107	0.00
52 CM	Toluene	50.000	58.295	-16.6#	109	0.00
53 T	t-1,3-Dichloropropene	50.000	57.105	-14.2	103	0.00
54 T	cis-1,3-Dichloropropene	50.000	57.561	-15.1	104	0.00
55 T	1,1,2-Trichloroethane	50.000	57.390	-14.8	108	0.00
56 T	Ethyl methacrylate	50.000	61.911	-23.8	105	0.00
57 T	1,3-Dichloropropane	50.000	57.876	-15.8	107	0.00
58 T	2-Chloroethyl Vinyl ether	250.000	304.242	-21.7	101	0.00
59 T	2-Hexanone	250.000	303.090	-21.2	105	0.00
60 T	Dibromochloromethane	50.000	59.046	-18.1	109	0.00
61 T	1,2-Dibromoethane	50.000	57.164	-14.3	104	0.00
62 S	4-Bromofluorobenzene	50.000	57.262	-14.5	107	0.00
63 I	Chlorobenzene-d5	50.000	50.000	0.0	101	0.00
64 T	Tetrachloroethene	50.000	45.415	9.2	97	0.00
65 PM	Chlorobenzene	50.000	52.142	-4.3	107	0.00
66 T	1,1,1,2-Tetrachloroethane	50.000	54.384	-8.8	109	0.00
67 C	Ethyl Benzene	50.000	53.794	-7.6#	110	0.00
68 T	m/p-Xylenes	100.000	110.067	-10.1	111	0.00
69 T	o-Xylene	50.000	53.371	-6.7	108	0.00
70 T	Styrene	50.000	57.477	-15.0	110	0.00
71 P	Bromoform	50.000	56.405	-12.8	111	0.00
72 I	1,4-Dichlorobenzene-d4	50.000	50.000	0.0	101	0.00
73 T	Isopropylbenzene	50.000	54.424	-8.8	112	0.00
74 T	N-ethyl acetate	50.000	40.223	19.6	98	0.00
75 P	1,1,2,2-Tetrachloroethane	50.000	55.070	-10.1	118	0.00
76 T	1,2,3-Trichloropropane	50.000	53.717	-7.4	106	0.00
77 T	Bromobenzene	50.000	55.390	-10.8	109	0.00
78 T	n-propylbenzene	50.000	54.998	-10.0	113	0.00
79 T	2-Chlorotoluene	50.000	52.964	-5.9	109	0.00
80 T	1,3,5-Trimethylbenzene	50.000	55.887	-11.8	113	0.00
81 T	trans-1,4-Dichloro-2-butene	50.000	57.865	-15.7	112	0.00
82 T	4-Chlorotoluene	50.000	56.933	-13.9	110	0.00
83 T	tert-Butylbenzene	50.000	53.710	-7.4	112	0.00
84 T	1,2,4-Trimethylbenzene	50.000	57.007	-14.0	112	0.00
85 T	sec-Butylbenzene	50.000	53.872	-7.7	113	0.00
86 T	p-Isopropyltoluene	50.000	54.723	-9.4	111	0.00
87 T	1,3-Dichlorobenzene	50.000	54.348	-8.7	112	0.00
88 T	1,4-Dichlorobenzene	50.000	52.479	-5.0	109	0.00
89 T	n-Butylbenzene	50.000	53.590	-7.2	112	0.00
90 T	Hexachloroethane	50.000	54.130	-8.3	114	0.00
91 T	1,2-Dichlorobenzene	50.000	53.690	-7.4	110	0.00
92 T	1,2-Dibromo-3-Chloropropane	50.000	52.428	-4.9	106	0.00
93 T	1,2,4-Trichlorobenzene	50.000	50.142	-0.3	105	0.00
94 T	Hexachlorobutadiene	50.000	51.883	-3.8	117	0.00
95 T	Naphthalene	50.000	53.119	-6.2	104	0.00

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN120425\
Data File : VN088437.D
Acq On : 04 Dec 2025 11:44
Operator : JC\MD
Sample : VSTDCCC050
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 3 Sample Multiplier: 1

Instrument :
MSVOA_N
LabSampleId :
VSTDCCC050

Quant Time: Dec 05 00:11:19 2025
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N112425W.M
Quant Title : SW846 8260
QLast Update : Tue Nov 25 01:34:52 2025
Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
Max. RRF Dev : 25% Max. Rel. Area : 150%

	Compound	Amount	Calc.	%Dev	Area%	Dev(min)
96 T	1,2,3-Trichlorobenzene	50.000	51.087	-2.2	105	0.00

(#) = Out of Range SPCC's out = 0 CCC's out = 6



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

VOLATILE CONTINUING CALIBRATION CHECK

Lab Name:	<u>Alliance</u>	Contract:	<u>JAC005</u>
Lab Code:	<u>ACE</u>	SDG No.:	<u>Q3757</u>
Instrument ID:	<u>MSVOA_X</u>	Calibration Date/Time:	<u>12/05/2025 08:45</u>
Lab File ID:	<u>VX048723.D</u>	Init. Calib. Date(s):	<u>12/04/2025 12/04/2025</u>
Heated Purge: (Y/N)	<u>N</u>	Init. Calib. Time(s):	<u>08:47 10:38</u>
GC Column:	<u>DB-624UI</u>	ID:	<u>0.18</u> (mm)

COMPOUND	RRF	RRF050	MIN RRF	%D	MAX%D
Vinyl Chloride	0.685	0.781		14.02	20
1,1-Dichloroethene	0.581	0.641		10.33	20
1,1-Dichloroethane	1.016	1.170	0.1	15.16	20
cis-1,2-Dichloroethene	0.669	0.724		8.22	20
1,1,1-Trichloroethane	0.975	0.958		-1.74	20
Benzene	1.398	1.412		1	20
1,2-Dichloroethane	0.495	0.470		-5.05	20
Trichloroethene	0.365	0.386		5.75	20
1,1,2-Trichloroethane	0.339	0.340		0.29	20
Tetrachloroethene	0.364	0.370		1.65	20
1,2-Dichloroethane-d4	0.671	0.588		-12.37	20
Dibromofluoromethane	0.344	0.344		0	20
Toluene-d8	1.207	1.162		-3.73	20
4-Bromofluorobenzene	0.416	0.430		3.37	20

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_X\Data\VX120525\
 Data File : VX048723.D
 Acq On : 05 Dec 2025 08:45
 Operator : JC/MD
 Sample : VSTDCCC050
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 VSTDCCC050

Manual Integrations
 APPROVED

Reviewed By : John Carlone 12/08/2025
 Supervised By : Semsettin Yesilyurt 12/08/2025

Quant Time: Dec 05 21:57:33 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X120425W.M
 Quant Title : SW846 8260
 QLast Update : Fri Dec 05 03:27:34 2025
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.525	168	218600	50.000	ug/l	-0.01
34) 1,4-Difluorobenzene	6.739	114	333007	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.037	117	285113	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.000	152	141895	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.928	65	128597	43.830	ug/l	-0.01
Spiked Amount 50.000	Range 78 - 117		Recovery =	87.660%		
35) Dibromofluoromethane	5.361	113	114548	49.958	ug/l	-0.01
Spiked Amount 50.000	Range 75 - 124		Recovery =	99.920%		
50) Toluene-d8	8.628	98	387114	48.165	ug/l	0.00
Spiked Amount 50.000	Range 92 - 112		Recovery =	96.320%		
62) 4-Bromofluorobenzene	11.061	95	143327	51.715	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery =	103.420%		
Target Compounds					Qvalue	
2) Dichlorodifluoromethane	1.172	85	145674	65.404	ug/l	99
3) Chloromethane	1.300	50	159260	56.664	ug/l	100
4) Vinyl Chloride	1.380	62	170695	57.025	ug/l	99
5) Bromomethane	1.611	94	112783	60.970	ug/l	98
6) Chloroethane	1.697	64	99323	53.315	ug/l	96
7) Trichlorofluoromethane	1.898	101	247779	57.756	ug/l	94
8) Diethyl Ether	2.130	74	90045	52.586	ug/l	98
9) 1,1,2-Trichlorotrifluo...	2.337	101	143083	57.531	ug/l	99
10) Methyl Iodide	2.453	142	192766	54.443	ug/l	99
11) Tert butyl alcohol	2.934	59	128210	270.659	ug/l	99
12) 1,1-Dichloroethene	2.325	96	140210	55.236	ug/l	98
13) Acrolein	2.233	56	115513	252.157	ug/l	99
14) Allyl chloride	2.660	41	253145	59.399	ug/l	99
15) Acrylonitrile	3.050	53	420739	285.052	ug/l	99
16) Acetone	2.367	43	252696	224.979	ug/l	99
17) Carbon Disulfide	2.514	76	407978	55.345	ug/l	99
18) Methyl Acetate	2.690	43	180477	57.109	ug/l	99
19) Methyl tert-butyl Ether	3.099	73	461178	55.825	ug/l	99
20) Methylene Chloride	2.788	84	147541	52.553	ug/l	98
21) trans-1,2-Dichloroethene	3.087	96	141112	54.468	ug/l	97
22) Diisopropyl ether	3.739	45	471234	59.164	ug/l	89
23) Vinyl Acetate	3.702	43	2070737	309.410	ug/l	97
24) 1,1-Dichloroethane	3.599	63	255858	57.605	ug/l	99
25) 2-Butanone	4.525	43	438661	264.615	ug/l	96
26) 2,2-Dichloropropane	4.458	77	221602	56.387	ug/l	97
27) cis-1,2-Dichloroethene	4.471	96	158350	54.165	ug/l	94
28) Bromochloromethane	4.873	49	95984	44.090	ug/l	99
29) Tetrahydrofuran	4.971	42	279279	220.633	ug/l	100
30) Chloroform	5.068	83	227566	47.817	ug/l	96
31) Cyclohexane	5.458	56	193584	48.816	ug/l	97
32) 1,1,1-Trichloroethane	5.367	97	209445	49.117	ug/l	99
36) 1,1-Dichloropropene	5.672	75	154341	49.146	ug/l	99
37) Ethyl Acetate	4.684	43	209277	55.770	ug/l	99
38) Carbon Tetrachloride	5.653	117	183592	51.612	ug/l	95
39) Methylcyclohexane	7.360	83	201583	54.284	ug/l	98
40) Benzene	6.013	78	470247	50.512	ug/l	100

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120525\
 Data File : VX048723.D
 Acq On : 05 Dec 2025 08:45
 Operator : JC/MD
 Sample : VSTDCCC050
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 VSTDCCC050

Manual Integrations
 APPROVED

Reviewed By :John Carlone 12/08/2025
 Supervised By :Semsettin Yesilyurt 12/08/2025

Quant Time: Dec 05 21:57:33 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X120425W.M
 Quant Title : SW846 8260
 QLast Update : Fri Dec 05 03:27:34 2025
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Methacrylonitrile	4.885	41	85389	50.012	ug/l	98
42) 1,2-Dichloroethane	6.062	62	156454	47.481	ug/l	99
43) Isopropyl Acetate	6.312	43	272582	47.972	ug/l	100
44) Trichloroethene	7.104	130	128510	52.904	ug/l	98
45) 1,2-Dichloropropane	7.409	63	113587	48.816	ug/l	98
46) Dibromomethane	7.562	93	90217	51.127	ug/l	98
47) Bromodichloromethane	7.799	83	174628	47.467	ug/l	98
48) Methyl methacrylate	7.671	41	132257	46.271	ug/l	96
49) 1,4-Dioxane	7.635	88	38976	991.484	ug/l	94
51) 4-Methyl-2-Pentanone	8.549	43	932226	264.320	ug/l	99
52) Toluene	8.695	92	288281	51.352	ug/l	100
53) t-1,3-Dichloropropene	8.958	75	177909	50.831	ug/l	99
54) cis-1,3-Dichloropropene	8.348	75	214885	57.128	ug/l	98
55) 1,1,2-Trichloroethane	9.134	97	113331	50.174	ug/l	97
56) Ethyl methacrylate	9.098	69	186000	52.609	ug/l	95
57) 1,3-Dichloropropane	9.287	76	185763	49.459	ug/l	100
58) 2-Chloroethyl Vinyl ether	8.220	63	541236	283.237	ug/l	100
59) 2-Hexanone	9.409	43	597075	241.164	ug/l	98
60) Dibromochloromethane	9.500	129	142497	51.018	ug/l	99
61) 1,2-Dibromoethane	9.592	107	123509	50.522	ug/l	99
64) Tetrachloroethene	9.256	164	105510	50.770	ug/l	97
65) Chlorobenzene	10.061	112	314975	50.470	ug/l	96
66) 1,1,1,2-Tetrachloroethane	10.146	131	116117	52.020	ug/l	99
67) Ethyl Benzene	10.177	91	546930	53.707	ug/l	100
68) m/p-Xylenes	10.281	106	420696	106.052	ug/l	98
69) o-Xylene	10.622	106	217816	58.597	ug/l	98
70) Styrene	10.634	104	372009	58.297	ug/l	99
71) Bromoform	10.780	173	104690	52.533	ug/l #	100
73) Isopropylbenzene	10.945	105	557420	58.295	ug/l	100
74) N-amyl acetate	10.823	43	263606	63.447	ug/l	99
75) 1,1,2,2-Tetrachloroethane	11.195	83	179135	55.522	ug/l	99
76) 1,2,3-Trichloropropane	11.219	75	149256m	56.476	ug/l	
77) Bromobenzene	11.177	156	138509	54.271	ug/l	99
78) n-propylbenzene	11.286	91	664210	59.976	ug/l	99
79) 2-Chlorotoluene	11.347	91	390949	57.891	ug/l	99
80) 1,3,5-Trimethylbenzene	11.433	105	424884	54.057	ug/l	99
81) trans-1,4-Dichloro-2-b...	11.000	75	63823	57.364	ug/l #	82
82) 4-Chlorotoluene	11.433	91	418818	53.145	ug/l	98
83) tert-Butylbenzene	11.695	119	442145	54.139	ug/l	99
84) 1,2,4-Trimethylbenzene	11.732	105	429499	54.171	ug/l	99
85) sec-Butylbenzene	11.872	105	543871	54.950	ug/l	98
86) p-Isopropyltoluene	11.988	119	464335	54.884	ug/l	99
87) 1,3-Dichlorobenzene	11.951	146	246588	52.411	ug/l	99
88) 1,4-Dichlorobenzene	12.018	146	246394	50.678	ug/l	99
89) n-Butylbenzene	12.311	91	421225	53.176	ug/l	99
90) Hexachloroethane	12.518	117	87435	50.636	ug/l	94
91) 1,2-Dichlorobenzene	12.317	146	233490	51.188	ug/l	98
92) 1,2-Dibromo-3-Chloropr...	12.920	75	36967	47.079	ug/l	93
93) 1,2,4-Trichlorobenzene	13.567	180	159178	52.838	ug/l	99
94) Hexachlorobutadiene	13.707	225	66243	53.159	ug/l	97
95) Naphthalene	13.756	128	491675	56.310	ug/l	100
96) 1,2,3-Trichlorobenzene	13.938	180	146739	52.816	ug/l	99

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120525\
Data File : VX048723.D
Acq On : 05 Dec 2025 08:45
Operator : JC/MD
Sample : VSTDCCC050
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 2 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
VSTDCCC050

Manual Integrations
APPROVED

Reviewed By :John Carlone 12/08/2025
Supervised By :Semsettin Yesilyurt 12/08/2025

Quant Time: Dec 05 21:57:33 2025
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X120425W.M
Quant Title : SW846 8260
QLast Update : Fri Dec 05 03:27:34 2025
Response via : Initial Calibration

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

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

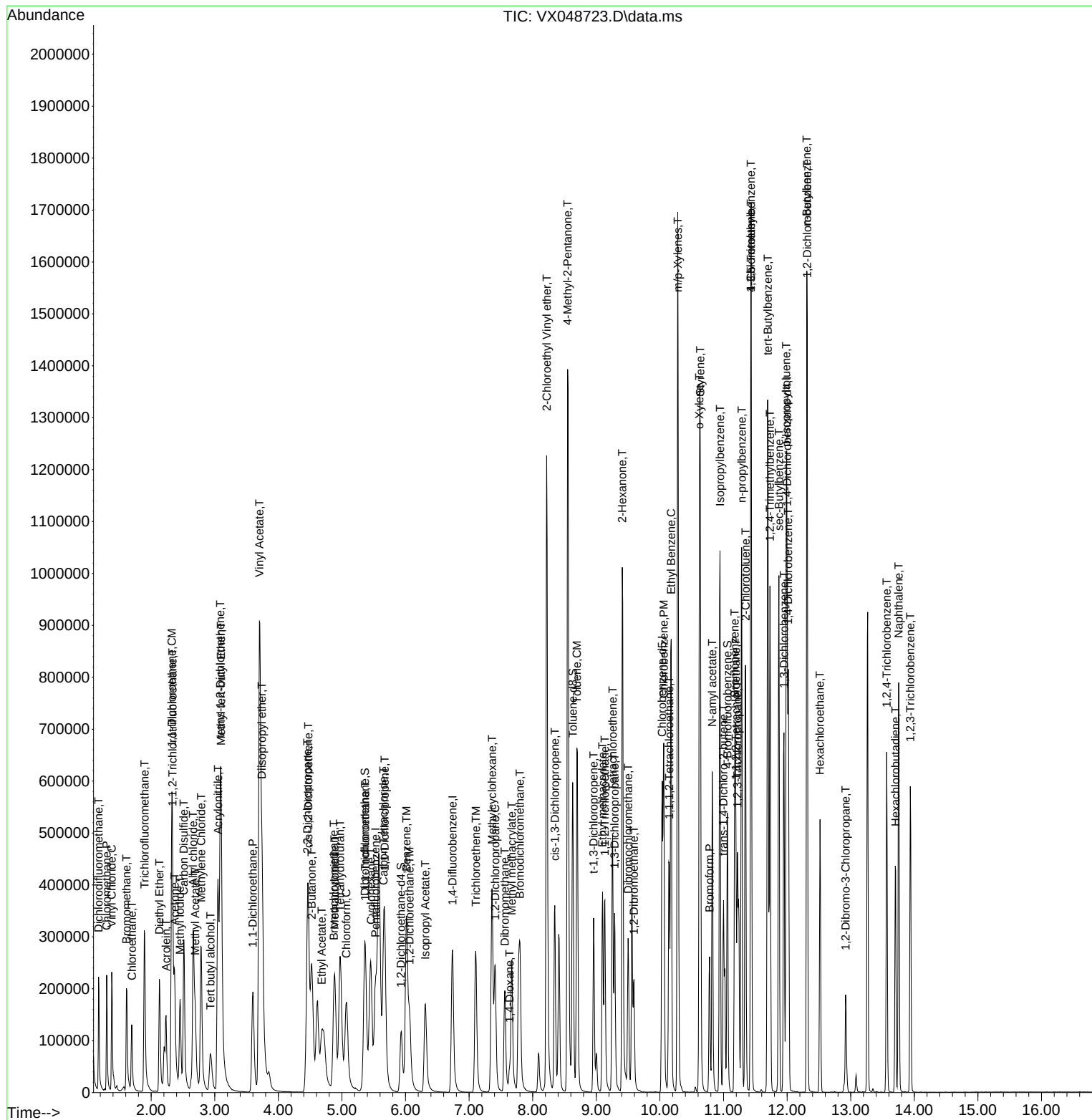
Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120525\
Data File : VX048723.D
Acq On : 05 Dec 2025 08:45
Operator : JC/MD
Sample : VSTDCCC050
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 2 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
VSTDCCC050

Manual Integrations APPROVED

Reviewed By :John Carlone 12/08/2025
Supervised By :Semsettin Yesilyurt 12/08/2025

Quant Time: Dec 05 21:57:33 2025
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X120425W.M
Quant Title : SW846 8260
QLast Update : Fri Dec 05 03:27:34 2025
Response via : Initial Calibration



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120525\
 Data File : VX048723.D
 Acq On : 05 Dec 2025 08:45
 Operator : JC/MD
 Sample : VSTDCCC050
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 MSVOA_X
 LabSampleId :
 VSTDCCC050

Quant Time: Dec 05 21:57:33 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X120425W.M
 Quant Title : SW846 8260
 QLast Update : Fri Dec 05 03:27:34 2025
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
 Max. RRF Dev : 25% Max. Rel. Area : 150%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
1 I	Pentafluorobenzene	1.000	1.000	0.0	117	-0.01
2 T	Dichlorodifluoromethane	0.509	0.666	-30.8#	137	0.00
3 P	Chloromethane	0.643	0.729	-13.4	132	0.00
4 C	Vinyl Chloride	0.685	0.781	-14.0#	129	0.00
5 T	Bromomethane	0.423	0.516	-22.0	129	0.00
6 T	Chloroethane	0.426	0.454	-6.6	125	0.00
7 T	Trichlorofluoromethane	0.981	1.133	-15.5	129	0.00
8 T	Diethyl Ether	0.392	0.412	-5.1	127	0.00
9 T	1,1,2-Trichlorotrifluoroeth	0.569	0.655	-15.1	129	0.00
10 T	Methyl Iodide	0.810	0.882	-8.9	119	0.00
11 T	Tert butyl alcohol	0.108	0.117	-8.3	120	0.00
12 CM	1,1-Dichloroethene	0.581	0.641	-10.3#	125	0.00
13 T	Acrolein	0.091	0.106	-16.5	144	0.00
14 T	Allyl chloride	0.975	1.158	-18.8	126	0.00
15 T	Acrylonitrile	0.338	0.385	-13.9	123	0.00
16 T	Acetone	0.257	0.231	10.1	110	0.00
17 T	Carbon Disulfide	1.686	1.866	-10.7	123	0.00
18 T	Methyl Acetate	0.723	0.826	-14.2	124	0.00
19 T	Methyl tert-butyl Ether	1.890	2.110	-11.6	124	0.00
20 T	Methylene Chloride	0.642	0.675	-5.1	122	0.00
21 T	trans-1,2-Dichloroethene	0.593	0.646	-8.9	121	0.00
22 T	Diisopropyl ether	1.822	2.156	-18.3	161#	-0.01
23 T	Vinyl Acetate	1.531	1.895	-23.8	161#	-0.01
24 P	1,1-Dichloroethane	1.016	1.170	-15.2	151#	0.00
25 T	2-Butanone	0.379	0.401	-5.8	139	-0.01
26 T	2,2-Dichloropropane	0.899	1.014	-12.8	143	-0.01
27 T	cis-1,2-Dichloroethene	0.669	0.724	-8.2	135	-0.01
28 T	Bromochloromethane	0.498	0.439	11.8	123	-0.01
29 T	Tetrahydrofuran	0.290	0.256	11.7	121	-0.01
30 C	Chloroform	1.089	1.041	4.4#	124	-0.01
31 T	Cyclohexane	0.907	0.886	2.3	130	0.00
32 T	1,1,1-Trichloroethane	0.975	0.958	1.7	123	0.00
33 S	1,2-Dichloroethane-d4	0.671	0.588	12.4	119	-0.01
34 I	1,4-Difluorobenzene	1.000	1.000	0.0	115	0.00
35 S	Dibromofluoromethane	0.344	0.344	0.0	119	-0.01
36 T	1,1-Dichloropropene	0.472	0.463	1.9	122	-0.01
37 T	Ethyl Acetate	0.563	0.628	-11.5	134	-0.01
38 T	Carbon Tetrachloride	0.534	0.551	-3.2	120	-0.01
39 T	Methylcyclohexane	0.558	0.605	-8.4	109	0.00
40 TM	Benzene	1.398	1.412	-1.0	123	-0.01
41 T	Methacrylonitrile	0.256	0.256	0.0	122	-0.02
42 TM	1,2-Dichloroethane	0.495	0.470	5.1	120	-0.01
43 T	Isopropyl Acetate	0.853	0.819	4.0	125	-0.01
44 TM	Trichloroethene	0.365	0.386	-5.8	121	0.00
45 C	1,2-Dichloropropane	0.349	0.341	2.3#	103	0.00
46 T	Dibromomethane	0.265	0.271	-2.3	108	0.00
47 T	Bromodichloromethane	0.552	0.524	5.1	103	0.00
48 T	Methyl methacrylate	0.429	0.397	7.5	98	0.00

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120525\
 Data File : VX048723.D
 Acq On : 05 Dec 2025 08:45
 Operator : JC/MD
 Sample : VSTDCCC050
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 MSVOA_X
 LabSampleId :
 VSTDCCC050

Quant Time: Dec 05 21:57:33 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X120425W.M
 Quant Title : SW846 8260
 QLast Update : Fri Dec 05 03:27:34 2025
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
 Max. RRF Dev : 25% Max. Rel. Area : 150%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
49 T	1,4-Dioxane	0.006	0.006	0.0	107	0.00
50 S	Toluene-d8	1.207	1.162	3.7	102	0.00
51 T	4-Methyl-2-Pentanone	0.530	0.560	-5.7	107	0.00
52 CM	Toluene	0.843	0.866	-2.7#	103	0.00
53 T	t-1,3-Dichloropropene	0.526	0.534	-1.5	103	0.00
54 T	cis-1,3-Dichloropropene	0.565	0.645	-14.2	115	0.00
55 T	1,1,2-Trichloroethane	0.339	0.340	-0.3	104	0.00
56 T	Ethyl methacrylate	0.531	0.559	-5.3	103	0.00
57 T	1,3-Dichloropropane	0.564	0.558	1.1	102	0.00
58 T	2-Chloroethyl Vinyl ether	0.287	0.325	-13.2	109	0.00
59 T	2-Hexanone	0.372	0.359	3.5	93	0.00
60 T	Dibromochloromethane	0.419	0.428	-2.1	105	0.00
61 T	1,2-Dibromoethane	0.367	0.371	-1.1	104	0.00
62 S	4-Bromofluorobenzene	0.416	0.430	-3.4	108	0.00
63 I	Chlorobenzene-d5	1.000	1.000	0.0	102	0.00
64 T	Tetrachloroethene	0.364	0.370	-1.6	107	0.00
65 PM	Chlorobenzene	1.094	1.105	-1.0	104	0.00
66 T	1,1,1,2-Tetrachloroethane	0.391	0.407	-4.1	110	0.00
67 C	Ethyl Benzene	1.786	1.918	-7.4#	104	0.00
68 T	m/p-Xylenes	0.696	0.738	-6.0	105	0.00
69 T	o-Xylene	0.652	0.764	-17.2	115	0.00
70 T	Styrene	1.119	1.305	-16.6	114	0.00
71 P	Bromoform	0.349	0.367	-5.2	114	0.00
72 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	104	0.00
73 T	Isopropylbenzene	3.369	3.928	-16.6	112	0.00
74 T	N-amyl acetate	1.464	1.858	-26.9#	111	0.00
75 P	1,1,2,2-Tetrachloroethane	1.137	1.262	-11.0	111	0.00
76 T	1,2,3-Trichloropropane	0.931	1.052	-13.0	114	0.00
77 T	Bromobenzene	0.899	0.976	-8.6	115	0.00
78 T	n-propylbenzene	3.902	4.681	-20.0	113	0.00
79 T	2-Chlorotoluene	2.380	2.755	-15.8	114	0.00
80 T	1,3,5-Trimethylbenzene	2.770	2.994	-8.1	106	0.00
81 T	trans-1,4-Dichloro-2-butene	0.392	0.450	-14.8	113	0.00
82 T	4-Chlorotoluene	2.777	2.952	-6.3	103	0.00
83 T	tert-Butylbenzene	2.878	3.116	-8.3	107	0.00
84 T	1,2,4-Trimethylbenzene	2.794	3.027	-8.3	105	0.00
85 T	sec-Butylbenzene	3.488	3.833	-9.9	106	0.00
86 T	p-Isopropyltoluene	2.981	3.272	-9.8	107	0.00
87 T	1,3-Dichlorobenzene	1.658	1.738	-4.8	110	0.00
88 T	1,4-Dichlorobenzene	1.713	1.736	-1.3	109	0.00
89 T	n-Butylbenzene	2.791	2.969	-6.4	104	0.00
90 T	Hexachloroethane	0.608	0.616	-1.3	108	0.00
91 T	1,2-Dichlorobenzene	1.607	1.646	-2.4	108	0.00
92 T	1,2-Dibromo-3-Chloropropane	0.277	0.261	5.8	101	0.00
93 T	1,2,4-Trichlorobenzene	1.062	1.122	-5.6	117	0.00
94 T	Hexachlorobutadiene	0.439	0.467	-6.4	123	0.00
95 T	Naphthalene	3.187	3.465	-8.7	115	0.00

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120525\
Data File : VX048723.D
Acq On : 05 Dec 2025 08:45
Operator : JC/MD
Sample : VSTDCCC050
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 2 Sample Multiplier: 1

Instrument :
MSVOA_X
LabSampleId :
VSTDCCC050

Quant Time: Dec 05 21:57:33 2025
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X120425W.M
Quant Title : SW846 8260
QLast Update : Fri Dec 05 03:27:34 2025
Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
Max. RRF Dev : 25% Max. Rel. Area : 150%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
96 T	1,2,3-Trichlorobenzene	0.979	1.034	-5.6	116	0.00

(#) = Out of Range

SPCC's out = 0 CCC's out = 6

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120525\
 Data File : VX048723.D
 Acq On : 05 Dec 2025 08:45
 Operator : JC/MD
 Sample : VSTDCCC050
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 MSVOA_X
 LabSampleId :
 VSTDCCC050

Quant Time: Dec 05 21:57:33 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X120425W.M
 Quant Title : SW846 8260
 QLast Update : Fri Dec 05 03:27:34 2025
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
 Max. RRF Dev : 25% Max. Rel. Area : 150%

	Compound	Amount	Calc.	%Dev	Area%	Dev(min)
1 I	Pentafluorobenzene	50.000	50.000	0.0	117	-0.01
2 T	Dichlorodifluoromethane	50.000	65.404	-30.8#	137	0.00
3 P	Chloromethane	50.000	56.664	-13.3	132	0.00
4 C	Vinyl Chloride	50.000	57.025	-14.0#	129	0.00
5 T	Bromomethane	50.000	60.970	-21.9	129	0.00
6 T	Chloroethane	50.000	53.315	-6.6	125	0.00
7 T	Trichlorofluoromethane	50.000	57.756	-15.5	129	0.00
8 T	Diethyl Ether	50.000	52.586	-5.2	127	0.00
9 T	1,1,2-Trichlorotrifluoroeth	50.000	57.531	-15.1	129	0.00
10 T	Methyl Iodide	50.000	54.443	-8.9	119	0.00
11 T	Tert butyl alcohol	250.000	270.659	-8.3	120	0.00
12 CM	1,1-Dichloroethene	50.000	55.236	-10.5#	125	0.00
13 T	Acrolein	250.000	252.157	-0.9	144	0.00
14 T	Allyl chloride	50.000	59.399	-18.8	126	0.00
15 T	Acrylonitrile	250.000	285.052	-14.0	123	0.00
16 T	Acetone	250.000	224.979	10.0	110	0.00
17 T	Carbon Disulfide	50.000	55.345	-10.7	123	0.00
18 T	Methyl Acetate	50.000	57.109	-14.2	124	0.00
19 T	Methyl tert-butyl Ether	50.000	55.825	-11.7	124	0.00
20 T	Methylene Chloride	50.000	52.553	-5.1	122	0.00
21 T	trans-1,2-Dichloroethene	50.000	54.468	-8.9	121	0.00
22 T	Diisopropyl ether	50.000	59.164	-18.3	161	-0.01
23 T	Vinyl Acetate	250.000	309.410	-23.8	161	-0.01
24 P	1,1-Dichloroethane	50.000	57.605	-15.2	151	0.00
25 T	2-Butanone	250.000	264.615	-5.8	139	-0.01
26 T	2,2-Dichloropropane	50.000	56.387	-12.8	143	-0.01
27 T	cis-1,2-Dichloroethene	50.000	54.165	-8.3	135	-0.01
28 T	Bromochloromethane	50.000	44.090	11.8	123	-0.01
29 T	Tetrahydrofuran	250.000	220.633	11.7	121	-0.01
30 C	Chloroform	50.000	47.817	4.4#	124	-0.01
31 T	Cyclohexane	50.000	48.816	2.4	130	0.00
32 T	1,1,1-Trichloroethane	50.000	49.117	1.8	123	0.00
33 S	1,2-Dichloroethane-d4	50.000	43.830	12.3	119	-0.01
34 I	1,4-Difluorobenzene	50.000	50.000	0.0	115	0.00
35 S	Dibromofluoromethane	50.000	49.958	0.1	119	-0.01
36 T	1,1-Dichloropropene	50.000	49.146	1.7	122	-0.01
37 T	Ethyl Acetate	50.000	55.770	-11.5	134	-0.01
38 T	Carbon Tetrachloride	50.000	51.612	-3.2	120	-0.01
39 T	Methylcyclohexane	50.000	54.284	-8.6	109	0.00
40 TM	Benzene	50.000	50.512	-1.0	123	-0.01
41 T	Methacrylonitrile	50.000	50.012	-0.0	122	-0.02
42 TM	1,2-Dichloroethane	50.000	47.481	5.0	120	-0.01
43 T	Isopropyl Acetate	50.000	47.972	4.1	125	-0.01
44 TM	Trichloroethene	50.000	52.904	-5.8	121	0.00
45 C	1,2-Dichloropropane	50.000	48.816	2.4#	103	0.00
46 T	Dibromomethane	50.000	51.127	-2.3	108	0.00
47 T	Bromodichloromethane	50.000	47.467	5.1	103	0.00
48 T	Methyl methacrylate	50.000	46.271	7.5	98	0.00

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120525\
 Data File : VX048723.D
 Acq On : 05 Dec 2025 08:45
 Operator : JC/MD
 Sample : VSTDCCC050
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 MSVOA_X
 LabSampleId :
 VSTDCCC050

Quant Time: Dec 05 21:57:33 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X120425W.M
 Quant Title : SW846 8260
 QLast Update : Fri Dec 05 03:27:34 2025
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
 Max. RRF Dev : 25% Max. Rel. Area : 150%

	Compound	Amount	Calc.	%Dev	Area%	Dev(min)
49 T	1,4-Dioxane	1000.000	991.484	0.9	107	0.00
50 S	Toluene-d8	50.000	48.165	3.7	102	0.00
51 T	4-Methyl-2-Pentanone	250.000	264.320	-5.7	107	0.00
52 CM	Toluene	50.000	51.352	-2.7#	103	0.00
53 T	t-1,3-Dichloropropene	50.000	50.831	-1.7	103	0.00
54 T	cis-1,3-Dichloropropene	50.000	57.128	-14.3	115	0.00
55 T	1,1,2-Trichloroethane	50.000	50.174	-0.3	104	0.00
56 T	Ethyl methacrylate	50.000	52.609	-5.2	103	0.00
57 T	1,3-Dichloropropane	50.000	49.459	1.1	102	0.00
58 T	2-Chloroethyl Vinyl ether	250.000	283.237	-13.3	109	0.00
59 T	2-Hexanone	250.000	241.164	3.5	93	0.00
60 T	Dibromochloromethane	50.000	51.018	-2.0	105	0.00
61 T	1,2-Dibromoethane	50.000	50.522	-1.0	104	0.00
62 S	4-Bromofluorobenzene	50.000	51.715	-3.4	108	0.00
63 I	Chlorobenzene-d5	50.000	50.000	0.0	102	0.00
64 T	Tetrachloroethene	50.000	50.770	-1.5	107	0.00
65 PM	Chlorobenzene	50.000	50.470	-0.9	104	0.00
66 T	1,1,1,2-Tetrachloroethane	50.000	52.020	-4.0	110	0.00
67 C	Ethyl Benzene	50.000	53.707	-7.4#	104	0.00
68 T	m/p-Xylenes	100.000	106.052	-6.1	105	0.00
69 T	o-Xylene	50.000	58.597	-17.2	115	0.00
70 T	Styrene	50.000	58.297	-16.6	114	0.00
71 P	Bromoform	50.000	52.533	-5.1	114	0.00
72 I	1,4-Dichlorobenzene-d4	50.000	50.000	0.0	104	0.00
73 T	Isopropylbenzene	50.000	58.295	-16.6	112	0.00
74 T	N-amyl acetate	50.000	63.447	-26.9#	111	0.00
75 P	1,1,2,2-Tetrachloroethane	50.000	55.522	-11.0	111	0.00
76 T	1,2,3-Trichloropropane	50.000	56.476	-13.0	114	0.00
77 T	Bromobenzene	50.000	54.271	-8.5	115	0.00
78 T	n-propylbenzene	50.000	59.976	-20.0	113	0.00
79 T	2-Chlorotoluene	50.000	57.891	-15.8	114	0.00
80 T	1,3,5-Trimethylbenzene	50.000	54.057	-8.1	106	0.00
81 T	trans-1,4-Dichloro-2-butene	50.000	57.364	-14.7	113	0.00
82 T	4-Chlorotoluene	50.000	53.145	-6.3	103	0.00
83 T	tert-Butylbenzene	50.000	54.139	-8.3	107	0.00
84 T	1,2,4-Trimethylbenzene	50.000	54.171	-8.3	105	0.00
85 T	sec-Butylbenzene	50.000	54.950	-9.9	106	0.00
86 T	p-Isopropyltoluene	50.000	54.884	-9.8	107	0.00
87 T	1,3-Dichlorobenzene	50.000	52.411	-4.8	110	0.00
88 T	1,4-Dichlorobenzene	50.000	50.678	-1.4	109	0.00
89 T	n-Butylbenzene	50.000	53.176	-6.4	104	0.00
90 T	Hexachloroethane	50.000	50.636	-1.3	108	0.00
91 T	1,2-Dichlorobenzene	50.000	51.188	-2.4	108	0.00
92 T	1,2-Dibromo-3-Chloropropane	50.000	47.079	5.8	101	0.00
93 T	1,2,4-Trichlorobenzene	50.000	52.838	-5.7	117	0.00
94 T	Hexachlorobutadiene	50.000	53.159	-6.3	123	0.00
95 T	Naphthalene	50.000	56.310	-12.6	115	0.00

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120525\
Data File : VX048723.D
Acq On : 05 Dec 2025 08:45
Operator : JC/MD
Sample : VSTDCCC050
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 2 Sample Multiplier: 1

Instrument :
MSVOA_X
LabSampleId :
VSTDCCC050

Quant Time: Dec 05 21:57:33 2025
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X120425W.M
Quant Title : SW846 8260
QLast Update : Fri Dec 05 03:27:34 2025
Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
Max. RRF Dev : 25% Max. Rel. Area : 150%

	Compound	Amount	Calc.	%Dev	Area%	Dev(min)
96 T	1,2,3-Trichlorobenzene	50.000	52.816	-5.6	116	0.00

(#) = Out of Range SPCC's out = 0 CCC's out = 6



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

VOLATILE CONTINUING CALIBRATION CHECK

Lab Name:	<u>Alliance</u>	Contract:	<u>JAC005</u>
Lab Code:	<u>ACE</u>	SDG No.:	<u>Q3757</u>
Instrument ID:	<u>MSVOA_X</u>	Calibration Date/Time:	<u>12/08/2025</u> <u>09:20</u>
Lab File ID:	<u>VX048742.D</u>	Init. Calib. Date(s):	<u>12/04/2025</u> <u>12/04/2025</u>
Heated Purge: (Y/N)	<u>N</u>	Init. Calib. Time(s):	<u>08:47</u> <u>10:38</u>
GC Column:	<u>DB-624UI</u>	ID:	<u>0.18</u> (mm)

COMPOUND	RRF	RRF050	MIN RRF	%D	MAX%D
Vinyl Chloride	0.685	0.671		-2.04	20
1,1-Dichloroethene	0.581	0.547		-5.85	20
1,1-Dichloroethane	1.016	1.086	0.1	6.89	20
cis-1,2-Dichloroethene	0.669	0.701		4.78	20
1,1,1-Trichloroethane	0.975	0.976		0.1	20
Benzene	1.398	1.441		3.08	20
1,2-Dichloroethane	0.495	0.558		12.73	20
Trichloroethene	0.365	0.370		1.37	20
1,1,2-Trichloroethane	0.339	0.389		14.75	20
Tetrachloroethene	0.364	0.345		-5.22	20
1,2-Dichloroethane-d4	0.671	0.665		-0.89	20
Dibromofluoromethane	0.344	0.346		0.58	20
Toluene-d8	1.207	1.218		0.91	20
4-Bromofluorobenzene	0.416	0.450		8.17	20

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_X\Data\VX120825\
 Data File : VX048742.D
 Acq On : 08 Dec 2025 09:20
 Operator : JC/MD
 Sample : VSTDCCC050
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 VSTDCCC050

Manual Integrations
 APPROVED

Reviewed By :John Carlone 12/09/2025
 Supervised By :Mahesh Dadoda 12/09/2025

Quant Time: Dec 09 04:05:01 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X120425W.M
 Quant Title : SW846 8260
 QLast Update : Fri Dec 05 03:27:34 2025
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.531	168	174813	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.739	114	271670	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.037	117	259893	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.000	152	132353	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.934	65	116232	49.538	ug/l	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery = 99.080%			
35) Dibromofluoromethane	5.361	113	93905	50.202	ug/l	-0.01
Spiked Amount 50.000	Range 75 - 124		Recovery = 100.400%			
50) Toluene-d8	8.628	98	330988	50.480	ug/l	0.00
Spiked Amount 50.000	Range 92 - 112		Recovery = 100.960%			
62) 4-Bromofluorobenzene	11.061	95	122254	54.070	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery = 108.140%			
Target Compounds					Qvalue	
2) Dichlorodifluoromethane	1.172	85	109463	61.456	ug/l	98
3) Chloromethane	1.300	50	112158	49.901	ug/l	99
4) Vinyl Chloride	1.380	62	117239	48.978	ug/l	97
5) Bromomethane	1.611	94	79697	53.876	ug/l	97
6) Chloroethane	1.691	64	70486	47.313	ug/l	96
7) Trichlorofluoromethane	1.892	101	168328	49.064	ug/l	95
8) Diethyl Ether	2.130	74	68655	50.137	ug/l	99
9) 1,1,2-Trichlorotrifluo...	2.331	101	96741	48.641	ug/l	99
10) Methyl Iodide	2.453	142	142606	50.365	ug/l	98
11) Tert butyl alcohol	2.934	59	104607	276.145	ug/l	99
12) 1,1-Dichloroethene	2.319	96	95593	47.091	ug/l	98
13) Acrolein	2.233	56	122007	320.155	ug/l	100
14) Allyl chloride	2.660	41	182879	53.660	ug/l	96
15) Acrylonitrile	3.050	53	323941	274.443	ug/l	99
16) Acetone	2.367	43	266788	297.020	ug/l	97
17) Carbon Disulfide	2.514	76	276616	46.924	ug/l	99
18) Methyl Acetate	2.690	43	147186	58.240	ug/l	98
19) Methyl tert-butyl Ether	3.099	73	346867	52.505	ug/l	98
20) Methylene Chloride	2.782	84	106567	47.466	ug/l	97
21) trans-1,2-Dichloroethene	3.087	96	96842	46.743	ug/l	97
22) Diisopropyl ether	3.745	45	361202	56.708	ug/l	97
23) Vinyl Acetate	3.702	43	1644500	307.270	ug/l	97
24) 1,1-Dichloroethane	3.599	63	189815	53.440	ug/l	99
25) 2-Butanone	4.525	43	409699	309.048	ug/l	95
26) 2,2-Dichloropropane	4.452	77	163755	52.104	ug/l	99
27) cis-1,2-Dichloroethene	4.471	96	122467	52.383	ug/l	94
28) Bromochloromethane	4.879	49	84093	48.303	ug/l	89
29) Tetrahydrofuran	4.971	42	292738	289.193	ug/l	96
30) Chloroform	5.068	83	201485	52.941	ug/l	100
31) Cyclohexane	5.452	56	152479	48.082	ug/l	96
32) 1,1,1-Trichloroethane	5.361	97	170606	50.030	ug/l	97
36) 1,1-Dichloropropene	5.672	75	129381	50.500	ug/l	99
37) Ethyl Acetate	4.684	43	200103	65.365	ug/l	97
38) Carbon Tetrachloride	5.659	117	151224	52.111	ug/l	93
39) Methylcyclohexane	7.360	83	159338	52.596	ug/l	99
40) Benzene	6.013	78	391521	51.550	ug/l	98

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120825\
 Data File : VX048742.D
 Acq On : 08 Dec 2025 09:20
 Operator : JC/MD
 Sample : VSTDCCC050
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 2 Sample Multiplier: 1

Instrument :

MSVOA_X

ClientSampleId :

VSTDCCC050

Manual Integrations

APPROVED

Reviewed By :John Carlone 12/09/2025

Supervised By :Mahesh Dadoda 12/09/2025

Quant Time: Dec 09 04:05:01 2025

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X120425W.M

Quant Title : SW846 8260

QLast Update : Fri Dec 05 03:27:34 2025

Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Methacrylonitrile	4.891	41	87479	62.805	ug/l	97
42) 1,2-Dichloroethane	6.062	62	151479	56.351	ug/l	96
43) Isopropyl Acetate	6.312	43	269086	58.049	ug/l	96
44) Trichloroethene	7.104	130	100636	50.782	ug/l	93
45) 1,2-Dichloropropane	7.409	63	103675	54.615	ug/l	97
46) Dibromomethane	7.562	93	81979	56.948	ug/l	100
47) Bromodichloromethane	7.799	83	161338	53.756	ug/l	100
48) Methyl methacrylate	7.671	41	129988	55.745	ug/l	98
49) 1,4-Dioxane	7.641	88	41910	1306.826	ug/l	93
51) 4-Methyl-2-Pentanone	8.549	43	839339	291.714	ug/l	99
52) Toluene	8.702	92	259442	56.649	ug/l	100
53) t-1,3-Dichloropropene	8.958	75	166878	58.445	ug/l	99
54) cis-1,3-Dichloropropene	8.348	75	176260	57.439	ug/l	96
55) 1,1,2-Trichloroethane	9.134	97	105645	57.331	ug/l	100
56) Ethyl methacrylate	9.098	69	176042	61.034	ug/l	97
57) 1,3-Dichloropropane	9.287	76	173763	56.709	ug/l	100
58) 2-Chloroethyl Vinyl ether	8.220	63	502570	322.382	ug/l	99
59) 2-Hexanone	9.409	43	605996	300.030	ug/l	99
60) Dibromochloromethane	9.500	129	132496	58.148	ug/l	99
61) 1,2-Dibromoethane	9.592	107	115535	57.930	ug/l	100
64) Tetrachloroethene	9.256	164	89617	47.307	ug/l	97
65) Chlorobenzene	10.061	112	287596	50.555	ug/l	99
66) 1,1,1,2-Tetrachloroethane	10.140	131	103746	50.988	ug/l	99
67) Ethyl Benzene	10.177	91	476354	51.315	ug/l	99
68) m/p-Xylenes	10.281	106	368936	102.029	ug/l	98
69) o-Xylene	10.622	106	180357	53.228	ug/l	98
70) Styrene	10.634	104	312915	53.795	ug/l	99
71) Bromoform	10.780	173	97762	53.817	ug/l	99
73) Isopropylbenzene	10.945	105	456673	51.202	ug/l	99
74) N-amyl acetate	10.823	43	218050	56.266	ug/l	97
75) 1,1,2,2-Tetrachloroethane	11.195	83	161398	53.631	ug/l	99
76) 1,2,3-Trichloropropane	11.219	75	132368m	53.696	ug/l	
77) Bromobenzene	11.177	156	122590	51.496	ug/l	96
78) n-propylbenzene	11.286	91	530512	51.357	ug/l	98
79) 2-Chlorotoluene	11.347	91	320717	50.915	ug/l	97
80) 1,3,5-Trimethylbenzene	11.433	105	379802	51.805	ug/l	100
81) trans-1,4-Dichloro-2-b...	11.000	75	58364	56.240	ug/l #	54
82) 4-Chlorotoluene	11.433	91	380068	51.705	ug/l	98
83) tert-Butylbenzene	11.695	119	387442	50.861	ug/l	98
84) 1,2,4-Trimethylbenzene	11.732	105	386338	52.241	ug/l	98
85) sec-Butylbenzene	11.872	105	471593	51.082	ug/l	99
86) p-Isopropyltoluene	11.988	119	409455	51.886	ug/l	99
87) 1,3-Dichlorobenzene	11.951	146	224294	51.109	ug/l	100
88) 1,4-Dichlorobenzene	12.018	146	227243	50.109	ug/l	99
89) n-Butylbenzene	12.311	91	368851	49.922	ug/l	99
90) Hexachloroethane	12.518	117	77355	48.028	ug/l	96
91) 1,2-Dichlorobenzene	12.317	146	216556	50.899	ug/l	100
92) 1,2-Dibromo-3-Chloropr...	12.920	75	37796	51.605	ug/l	99
93) 1,2,4-Trichlorobenzene	13.567	180	144749	51.512	ug/l	99
94) Hexachlorobutadiene	13.707	225	55748	47.962	ug/l	96
95) Naphthalene	13.756	128	458027	56.250	ug/l	99
96) 1,2,3-Trichlorobenzene	13.938	180	140730	54.305	ug/l	97

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120825\
Data File : VX048742.D
Acq On : 08 Dec 2025 09:20
Operator : JC/MD
Sample : VSTDCCC050
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 2 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
VSTDCCC050

Manual Integrations
APPROVED

Reviewed By :John Carlone 12/09/2025
Supervised By :Mahesh Dadoda 12/09/2025

Quant Time: Dec 09 04:05:01 2025
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X120425W.M
Quant Title : SW846 8260
QLast Update : Fri Dec 05 03:27:34 2025
Response via : Initial Calibration

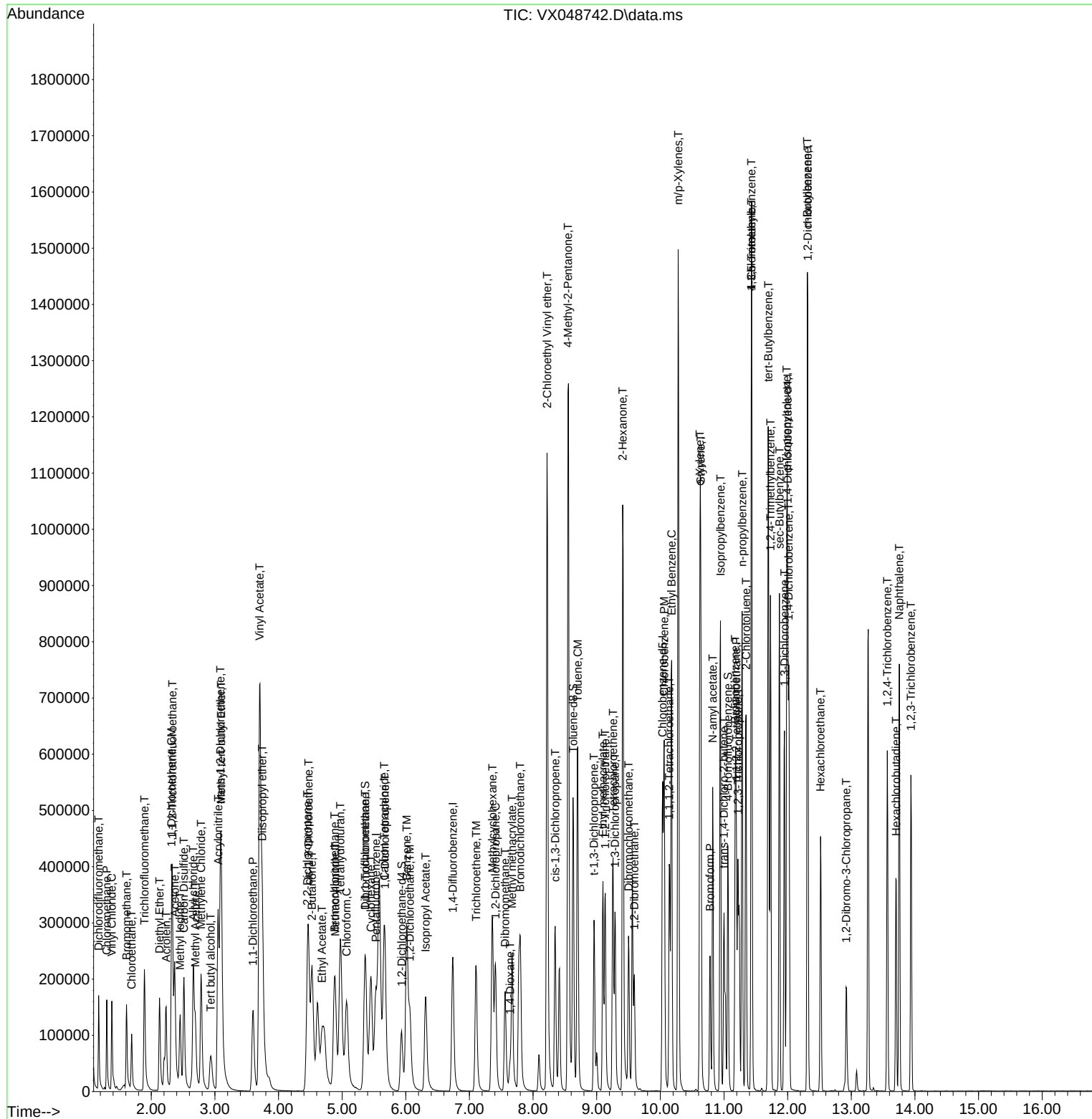
Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
----------	------	------	----------	------	-------	----------

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

Instrument :
MSVOA_X
ClientSampleId :
VSTDCCC050

Manual Integrations APPROVED

Reviewed By :John Carlone 12/09/2025
Supervised By :Mahesh Dadoda 12/09/2025



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120825\
 Data File : VX048742.D
 Acq On : 08 Dec 2025 09:20
 Operator : JC/MD
 Sample : VSTDCCC050
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 MSVOA_X
 LabSampleId :
 VSTDCCC050

Quant Time: Dec 09 04:05:01 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X120425W.M
 Quant Title : SW846 8260
 QLast Update : Fri Dec 05 03:27:34 2025
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
 Max. RRF Dev : 25% Max. Rel. Area : 150%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
1 I	Pentafluorobenzene	1.000	1.000	0.0	94	0.00
2 T	Dichlorodifluoromethane	0.509	0.626	-23.0	103	0.00
3 P	Chloromethane	0.643	0.642	0.2	93	0.00
4 C	Vinyl Chloride	0.685	0.671	2.0#	89	0.00
5 T	Bromomethane	0.423	0.456	-7.8	91	0.00
6 T	Chloroethane	0.426	0.403	5.4	89	0.00
7 T	Trichlorofluoromethane	0.981	0.963	1.8	87	0.00
8 T	Diethyl Ether	0.392	0.393	-0.3	97	0.00
9 T	1,1,2-Trichlorotrifluoroeth	0.569	0.553	2.8	87	0.00
10 T	Methyl Iodide	0.810	0.816	-0.7	88	0.00
11 T	Tert butyl alcohol	0.108	0.120	-11.1	98	0.00
12 CM	1,1-Dichloroethene	0.581	0.547	5.9#	85	0.00
13 T	Acrolein	0.091	0.140	-53.8#	153#	0.00
14 T	Allyl chloride	0.975	1.046	-7.3	91	0.00
15 T	Acrylonitrile	0.338	0.371	-9.8	95	0.00
16 T	Acetone	0.257	0.305	-18.7	116	0.00
17 T	Carbon Disulfide	1.686	1.582	6.2	84	0.00
18 T	Methyl Acetate	0.723	0.842	-16.5	101	0.00
19 T	Methyl tert-butyl Ether	1.890	1.984	-5.0	93	0.00
20 T	Methylene Chloride	0.642	0.610	5.0	88	0.00
21 T	trans-1,2-Dichloroethene	0.593	0.554	6.6	83	0.00
22 T	Diisopropyl ether	1.822	2.066	-13.4	124	0.00
23 T	Vinyl Acetate	1.531	1.881	-22.9	128	-0.01
24 P	1,1-Dichloroethane	1.016	1.086	-6.9	112	0.00
25 T	2-Butanone	0.379	0.469	-23.7	130	-0.01
26 T	2,2-Dichloropropane	0.899	0.937	-4.2	106	-0.02
27 T	cis-1,2-Dichloroethene	0.669	0.701	-4.8	104	-0.01
28 T	Bromochloromethane	0.498	0.481	3.4	108	0.00
29 T	Tetrahydrofuran	0.290	0.335	-15.5	127	-0.01
30 C	Chloroform	1.089	1.153	-5.9#	110	-0.01
31 T	Cyclohexane	0.907	0.872	3.9	103	-0.01
32 T	1,1,1-Trichloroethane	0.975	0.976	-0.1	100	0.00
33 S	1,2-Dichloroethane-d4	0.671	0.665	0.9	108	0.00
34 I	1,4-Difluorobenzene	1.000	1.000	0.0	94	0.00
35 S	Dibromofluoromethane	0.344	0.346	-0.6	97	-0.01
36 T	1,1-Dichloropropene	0.472	0.476	-0.8	102	-0.01
37 T	Ethyl Acetate	0.563	0.737	-30.9#	128	-0.01
38 T	Carbon Tetrachloride	0.534	0.557	-4.3	99	0.00
39 T	Methylcyclohexane	0.558	0.587	-5.2	86	0.00
40 TM	Benzene	1.398	1.441	-3.1	102	-0.01
41 T	Methacrylonitrile	0.256	0.322	-25.8#	125	-0.01
42 TM	1,2-Dichloroethane	0.495	0.558	-12.7	116	-0.01
43 T	Isopropyl Acetate	0.853	0.990	-16.1	123	-0.01
44 TM	Trichloroethene	0.365	0.370	-1.4	95	0.00
45 C	1,2-Dichloropropane	0.349	0.382	-9.5#	94	0.00
46 T	Dibromomethane	0.265	0.302	-14.0	98	0.00
47 T	Bromodichloromethane	0.552	0.594	-7.6	95	0.00
48 T	Methyl methacrylate	0.429	0.478	-11.4	97	0.00

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120825\
 Data File : VX048742.D
 Acq On : 08 Dec 2025 09:20
 Operator : JC/MD
 Sample : VSTDCCC050
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 MSVOA_X
 LabSampleId :
 VSTDCCC050

Quant Time: Dec 09 04:05:01 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X120425W.M
 Quant Title : SW846 8260
 QLast Update : Fri Dec 05 03:27:34 2025
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
 Max. RRF Dev : 25% Max. Rel. Area : 150%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
49 T	1,4-Dioxane	0.006	0.008	-33.3#	115	0.00
50 S	Toluene-d8	1.207	1.218	-0.9	87	0.00
51 T	4-Methyl-2-Pentanone	0.530	0.618	-16.6	96	0.00
52 CM	Toluene	0.843	0.955	-13.3#	93	0.00
53 T	t-1,3-Dichloropropene	0.526	0.614	-16.7	96	0.00
54 T	cis-1,3-Dichloropropene	0.565	0.649	-14.9	95	0.00
55 T	1,1,2-Trichloroethane	0.339	0.389	-14.7	97	0.00
56 T	Ethyl methacrylate	0.531	0.648	-22.0	97	0.00
57 T	1,3-Dichloropropane	0.564	0.640	-13.5	96	0.00
58 T	2-Chloroethyl Vinyl ether	0.287	0.370	-28.9#	101	0.00
59 T	2-Hexanone	0.372	0.446	-19.9	95	0.00
60 T	Dibromochloromethane	0.419	0.488	-16.5	98	0.00
61 T	1,2-Dibromoethane	0.367	0.425	-15.8	97	0.00
62 S	4-Bromofluorobenzene	0.416	0.450	-8.2	92	0.00
63 I	Chlorobenzene-d5	1.000	1.000	0.0	93	0.00
64 T	Tetrachloroethene	0.364	0.345	5.2	91	0.00
65 PM	Chlorobenzene	1.094	1.107	-1.2	95	0.00
66 T	1,1,1,2-Tetrachloroethane	0.391	0.399	-2.0	98	0.00
67 C	Ethyl Benzene	1.786	1.833	-2.6#	90	0.00
68 T	m/p-Xylenes	0.696	0.710	-2.0	92	0.00
69 T	o-Xylene	0.652	0.694	-6.4	95	0.00
70 T	Styrene	1.119	1.204	-7.6	96	0.00
71 P	Bromoform	0.349	0.376	-7.7	107	0.00
72 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	97	0.00
73 T	Isopropylbenzene	3.369	3.450	-2.4	92	0.00
74 T	N-ethyl acetate	1.464	1.647	-12.5	92	0.00
75 P	1,1,2,2-Tetrachloroethane	1.137	1.219	-7.2	100	0.00
76 T	1,2,3-Trichloropropane	0.931	1.000	-7.4	101	0.00
77 T	Bromobenzene	0.899	0.926	-3.0	101	0.00
78 T	n-propylbenzene	3.902	4.008	-2.7	90	0.00
79 T	2-Chlorotoluene	2.380	2.423	-1.8	93	0.00
80 T	1,3,5-Trimethylbenzene	2.770	2.870	-3.6	94	0.00
81 T	trans-1,4-Dichloro-2-butene	0.392	0.441	-12.5	103	0.00
82 T	4-Chlorotoluene	2.777	2.872	-3.4	93	0.00
83 T	tert-Butylbenzene	2.878	2.927	-1.7	94	0.00
84 T	1,2,4-Trimethylbenzene	2.794	2.919	-4.5	95	0.00
85 T	sec-Butylbenzene	3.488	3.563	-2.2	92	0.00
86 T	p-Isopropyltoluene	2.981	3.094	-3.8	94	0.00
87 T	1,3-Dichlorobenzene	1.658	1.695	-2.2	100	0.00
88 T	1,4-Dichlorobenzene	1.713	1.717	-0.2	100	0.00
89 T	n-Butylbenzene	2.791	2.787	0.1	91	0.00
90 T	Hexachloroethane	0.608	0.584	3.9	95	0.00
91 T	1,2-Dichlorobenzene	1.607	1.636	-1.8	100	0.00
92 T	1,2-Dibromo-3-Chloropropane	0.277	0.286	-3.2	103	0.00
93 T	1,2,4-Trichlorobenzene	1.062	1.094	-3.0	106	0.00
94 T	Hexachlorobutadiene	0.439	0.421	4.1	104	0.00
95 T	Naphthalene	3.187	3.461	-8.6	107	0.00

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120825\
Data File : VX048742.D
Acq On : 08 Dec 2025 09:20
Operator : JC/MD
Sample : VSTDCCC050
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 2 Sample Multiplier: 1

Instrument :
MSVOA_X
LabSampleId :
VSTDCCC050

Quant Time: Dec 09 04:05:01 2025
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X120425W.M
Quant Title : SW846 8260
QLast Update : Fri Dec 05 03:27:34 2025
Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
Max. RRF Dev : 25% Max. Rel. Area : 150%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
96 T	1,2,3-Trichlorobenzene	0.979	1.063	-8.6	111	0.00

(#) = Out of Range

SPCC's out = 0 CCC's out = 6

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120825\
 Data File : VX048742.D
 Acq On : 08 Dec 2025 09:20
 Operator : JC/MD
 Sample : VSTDCCC050
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 MSVOA_X
 LabSampleId :
 VSTDCCC050

Quant Time: Dec 09 04:05:01 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X120425W.M
 Quant Title : SW846 8260
 QLast Update : Fri Dec 05 03:27:34 2025
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
 Max. RRF Dev : 25% Max. Rel. Area : 150%

	Compound	Amount	Calc.	%Dev	Area%	Dev(min)
1 I	Pentafluorobenzene	50.000	50.000	0.0	94	0.00
2 T	Dichlorodifluoromethane	50.000	61.456	-22.9	103	0.00
3 P	Chloromethane	50.000	49.901	0.2	93	0.00
4 C	Vinyl Chloride	50.000	48.978	2.0#	89	0.00
5 T	Bromomethane	50.000	53.876	-7.8	91	0.00
6 T	Chloroethane	50.000	47.313	5.4	89	0.00
7 T	Trichlorofluoromethane	50.000	49.064	1.9	87	0.00
8 T	Diethyl Ether	50.000	50.137	-0.3	97	0.00
9 T	1,1,2-Trichlorotrifluoroeth	50.000	48.641	2.7	87	0.00
10 T	Methyl Iodide	50.000	50.365	-0.7	88	0.00
11 T	Tert butyl alcohol	250.000	276.145	-10.5	98	0.00
12 CM	1,1-Dichloroethene	50.000	47.091	5.8#	85	0.00
13 T	Acrolein	250.000	320.155	-28.1#	153	0.00
14 T	Allyl chloride	50.000	53.660	-7.3	91	0.00
15 T	Acrylonitrile	250.000	274.443	-9.8	95	0.00
16 T	Acetone	250.000	297.020	-18.8	116	0.00
17 T	Carbon Disulfide	50.000	46.924	6.2	84	0.00
18 T	Methyl Acetate	50.000	58.240	-16.5	101	0.00
19 T	Methyl tert-butyl Ether	50.000	52.505	-5.0	93	0.00
20 T	Methylene Chloride	50.000	47.466	5.1	88	0.00
21 T	trans-1,2-Dichloroethene	50.000	46.743	6.5	83	0.00
22 T	Diisopropyl ether	50.000	56.708	-13.4	124	0.00
23 T	Vinyl Acetate	250.000	307.270	-22.9	128	-0.01
24 P	1,1-Dichloroethane	50.000	53.440	-6.9	112	0.00
25 T	2-Butanone	250.000	309.048	-23.6	130	-0.01
26 T	2,2-Dichloropropane	50.000	52.104	-4.2	106	-0.02
27 T	cis-1,2-Dichloroethene	50.000	52.383	-4.8	104	-0.01
28 T	Bromochloromethane	50.000	48.303	3.4	108	0.00
29 T	Tetrahydrofuran	250.000	289.193	-15.7	127	-0.01
30 C	Chloroform	50.000	52.941	-5.9#	110	-0.01
31 T	Cyclohexane	50.000	48.082	3.8	103	-0.01
32 T	1,1,1-Trichloroethane	50.000	50.030	-0.1	100	0.00
33 S	1,2-Dichloroethane-d4	50.000	49.538	0.9	108	0.00
34 I	1,4-Difluorobenzene	50.000	50.000	0.0	94	0.00
35 S	Dibromofluoromethane	50.000	50.202	-0.4	97	-0.01
36 T	1,1-Dichloropropene	50.000	50.500	-1.0	102	-0.01
37 T	Ethyl Acetate	50.000	65.365	-30.7#	128	-0.01
38 T	Carbon Tetrachloride	50.000	52.111	-4.2	99	0.00
39 T	Methylcyclohexane	50.000	52.596	-5.2	86	0.00
40 TM	Benzene	50.000	51.550	-3.1	102	-0.01
41 T	Methacrylonitrile	50.000	62.805	-25.6#	125	-0.01
42 TM	1,2-Dichloroethane	50.000	56.351	-12.7	116	-0.01
43 T	Isopropyl Acetate	50.000	58.049	-16.1	123	-0.01
44 TM	Trichloroethene	50.000	50.782	-1.6	95	0.00
45 C	1,2-Dichloropropane	50.000	54.615	-9.2#	94	0.00
46 T	Dibromomethane	50.000	56.948	-13.9	98	0.00
47 T	Bromodichloromethane	50.000	53.756	-7.5	95	0.00
48 T	Methyl methacrylate	50.000	55.745	-11.5	97	0.00

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120825\
 Data File : VX048742.D
 Acq On : 08 Dec 2025 09:20
 Operator : JC/MD
 Sample : VSTDCCC050
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 MSVOA_X
 LabSampleId :
 VSTDCCC050

Quant Time: Dec 09 04:05:01 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X120425W.M
 Quant Title : SW846 8260
 QLast Update : Fri Dec 05 03:27:34 2025
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
 Max. RRF Dev : 25% Max. Rel. Area : 150%

	Compound	Amount	Calc.	%Dev	Area%	Dev(min)
49 T	1,4-Dioxane	1000.000	1306.826	-30.7#	115	0.00
50 S	Toluene-d8	50.000	50.480	-1.0	87	0.00
51 T	4-Methyl-2-Pentanone	250.000	291.714	-16.7	96	0.00
52 CM	Toluene	50.000	56.649	-13.3#	93	0.00
53 T	t-1,3-Dichloropropene	50.000	58.445	-16.9	96	0.00
54 T	cis-1,3-Dichloropropene	50.000	57.439	-14.9	95	0.00
55 T	1,1,2-Trichloroethane	50.000	57.331	-14.7	97	0.00
56 T	Ethyl methacrylate	50.000	61.034	-22.1	97	0.00
57 T	1,3-Dichloropropane	50.000	56.709	-13.4	96	0.00
58 T	2-Chloroethyl Vinyl ether	250.000	322.382	-29.0#	101	0.00
59 T	2-Hexanone	250.000	300.030	-20.0	95	0.00
60 T	Dibromochloromethane	50.000	58.148	-16.3	98	0.00
61 T	1,2-Dibromoethane	50.000	57.930	-15.9	97	0.00
62 S	4-Bromofluorobenzene	50.000	54.070	-8.1	92	0.00
63 I	Chlorobenzene-d5	50.000	50.000	0.0	93	0.00
64 T	Tetrachloroethene	50.000	47.307	5.4	91	0.00
65 PM	Chlorobenzene	50.000	50.555	-1.1	95	0.00
66 T	1,1,1,2-Tetrachloroethane	50.000	50.988	-2.0	98	0.00
67 C	Ethyl Benzene	50.000	51.315	-2.6#	90	0.00
68 T	m/p-Xylenes	100.000	102.029	-2.0	92	0.00
69 T	o-Xylene	50.000	53.228	-6.5	95	0.00
70 T	Styrene	50.000	53.795	-7.6	96	0.00
71 P	Bromoform	50.000	53.817	-7.6	107	0.00
72 I	1,4-Dichlorobenzene-d4	50.000	50.000	0.0	97	0.00
73 T	Isopropylbenzene	50.000	51.202	-2.4	92	0.00
74 T	N-amyl acetate	50.000	56.266	-12.5	92	0.00
75 P	1,1,2,2-Tetrachloroethane	50.000	53.631	-7.3	100	0.00
76 T	1,2,3-Trichloropropane	50.000	53.696	-7.4	101	0.00
77 T	Bromobenzene	50.000	51.496	-3.0	101	0.00
78 T	n-propylbenzene	50.000	51.357	-2.7	90	0.00
79 T	2-Chlorotoluene	50.000	50.915	-1.8	93	0.00
80 T	1,3,5-Trimethylbenzene	50.000	51.805	-3.6	94	0.00
81 T	trans-1,4-Dichloro-2-butene	50.000	56.240	-12.5	103	0.00
82 T	4-Chlorotoluene	50.000	51.705	-3.4	93	0.00
83 T	tert-Butylbenzene	50.000	50.861	-1.7	94	0.00
84 T	1,2,4-Trimethylbenzene	50.000	52.241	-4.5	95	0.00
85 T	sec-Butylbenzene	50.000	51.082	-2.2	92	0.00
86 T	p-Isopropyltoluene	50.000	51.886	-3.8	94	0.00
87 T	1,3-Dichlorobenzene	50.000	51.109	-2.2	100	0.00
88 T	1,4-Dichlorobenzene	50.000	50.109	-0.2	100	0.00
89 T	n-Butylbenzene	50.000	49.922	0.2	91	0.00
90 T	Hexachloroethane	50.000	48.028	3.9	95	0.00
91 T	1,2-Dichlorobenzene	50.000	50.899	-1.8	100	0.00
92 T	1,2-Dibromo-3-Chloropropane	50.000	51.605	-3.2	103	0.00
93 T	1,2,4-Trichlorobenzene	50.000	51.512	-3.0	106	0.00
94 T	Hexachlorobutadiene	50.000	47.962	4.1	104	0.00
95 T	Naphthalene	50.000	56.250	-12.5	107	0.00

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120825\
Data File : VX048742.D
Acq On : 08 Dec 2025 09:20
Operator : JC/MD
Sample : VSTDCCC050
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 2 Sample Multiplier: 1

Instrument :
MSVOA_X
LabSampleId :
VSTDCCC050

Quant Time: Dec 09 04:05:01 2025
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X120425W.M
Quant Title : SW846 8260
QLast Update : Fri Dec 05 03:27:34 2025
Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
Max. RRF Dev : 25% Max. Rel. Area : 150%

	Compound	Amount	Calc.	%Dev	Area%	Dev(min)
96 T	1,2,3-Trichlorobenzene	50.000	54.305	-8.6	111	0.00

(#) = Out of Range SPCC's out = 0 CCC's out = 6



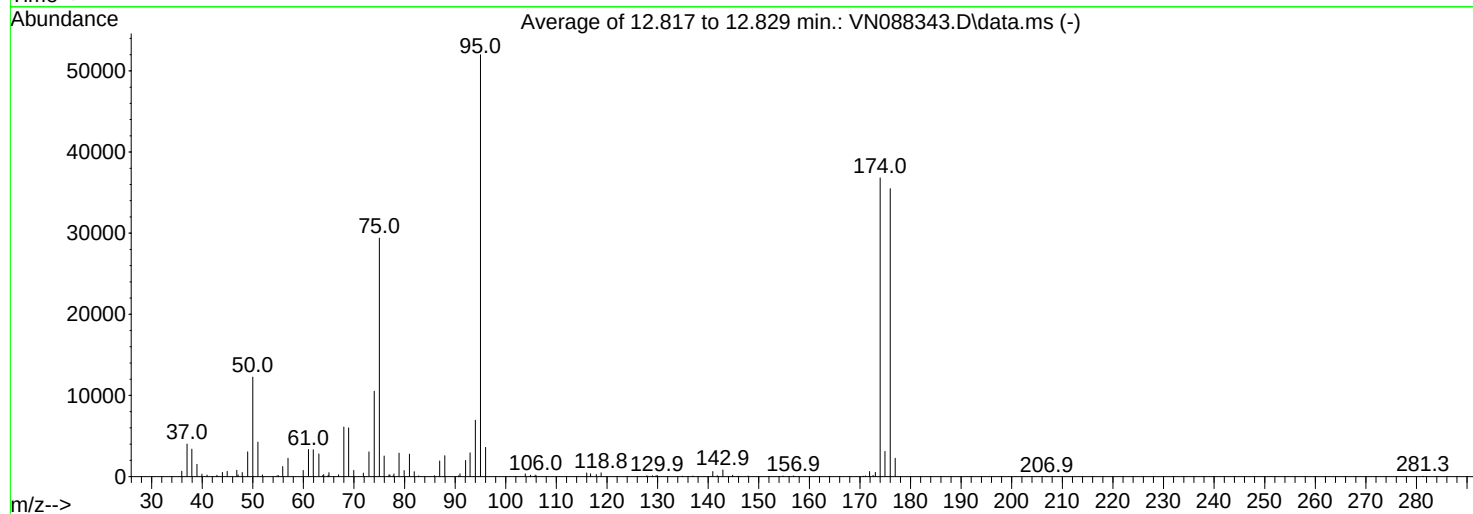
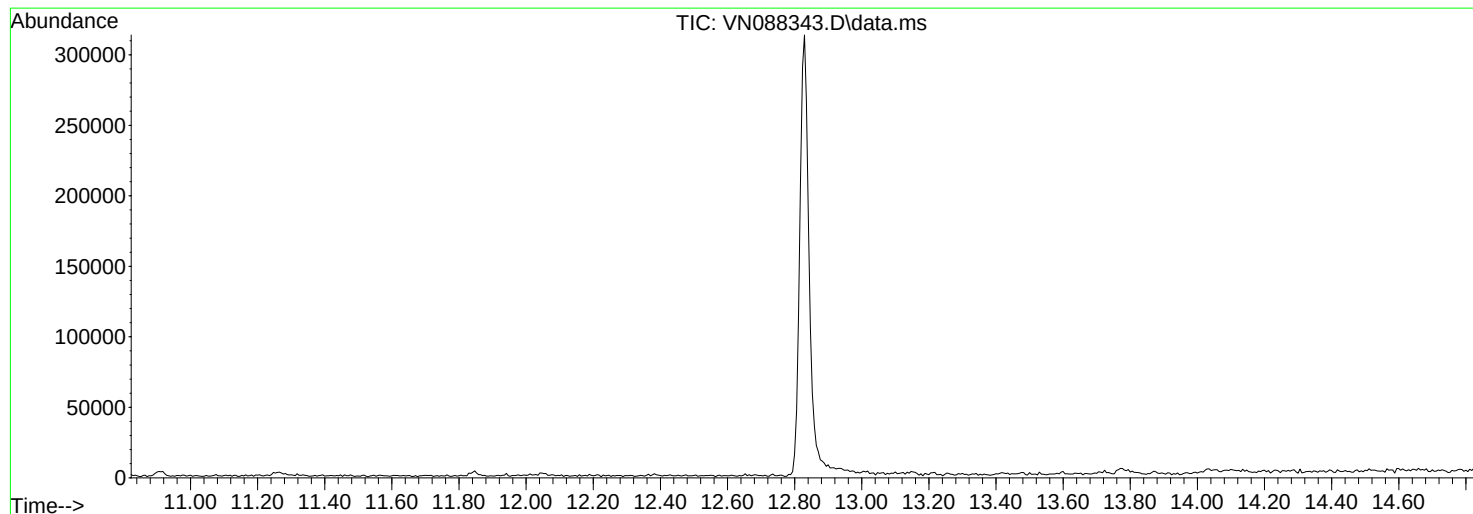
QC SAMPLE DATA

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN112425\
Data File : VN088343.D
Acq On : 24 Nov 2025 08:17
Operator : JC\MD
Sample : BFB
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
BFB

Integration File: RTEINT.P

Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N112425W.M
Title : SW846 8260
Last Update : Tue Nov 25 01:34:52 2025



AutoFind: Scans 1847, 1848, 1849; Background Corrected with Scan 1839

Target	Rel. to	Lower	Upper	Rel.	Raw	Result
Mass	Mass	Limit%	Limit%	Abn%	Abn	Pass/Fail
50	95	15	40	23.6	12249	PASS
75	95	30	60	56.6	29403	PASS
95	95	100	100	100.0	51987	PASS
96	95	5	9	6.9	3605	PASS
173	174	0.00	2	1.4	523	PASS
174	95	50	100	70.8	36808	PASS
175	174	5	9	8.4	3099	PASS
176	174	95	101	96.4	35493	PASS
177	176	5	9	6.4	2260	PASS

BFB

Modified:subtracted

m/z	abund.	m/z	abund.	m/z	abund.	m/z	abund.
35.95	660	47.90	508	61.95	3309	72.95	3051
37.00	4014	49.00	3039	63.05	2784	74.00	10530
37.95	3391	50.00	12249	63.80	113	75.00	29403
38.95	1502	51.00	4245	64.00	255	75.95	2531
39.95	311	51.90	213	64.90	158	76.90	202
40.95	180	54.80	110	65.05	474	77.10	198
42.90	161	55.10	103	66.95	229	77.70	63
44.00	531	55.90	1233	68.00	6110	77.90	320
44.95	637	56.95	2258	68.95	5997	78.90	2897
46.85	771	59.95	755	69.95	751	79.90	742
47.10	213	61.00	3348	71.85	421	80.95	2778

Average of 12.817 to 12.829 min.: VN088343.D\data.ms

BFB

Modified:subtracted

m/z	abund.	m/z	abund.	m/z	abund.	m/z	abund.
81.90	605	96.00	3605	129.70	84	156.90	87
82.80	89	103.85	320	129.90	147	170.70	51
85.90	119	104.90	161	134.90	52	171.10	82
86.95	1933	105.70	75	137.00	60	171.90	617
87.95	2571	105.95	190	140.90	629	172.60	144
90.70	103	116.00	445	141.80	113	173.05	523
90.95	348	116.75	347	142.90	817	174.00	36808
92.05	1981	117.90	250	144.85	179	174.95	3099
92.95	2935	118.85	465	145.90	56	176.00	35493
94.00	6957	127.90	108	147.90	68	176.95	2260
95.00	51987	129.00	114	154.80	69	206.95	15

Average of 12.817 to 12.829 min.: VN088343.D\data.ms

BFB

Modified:subtracted

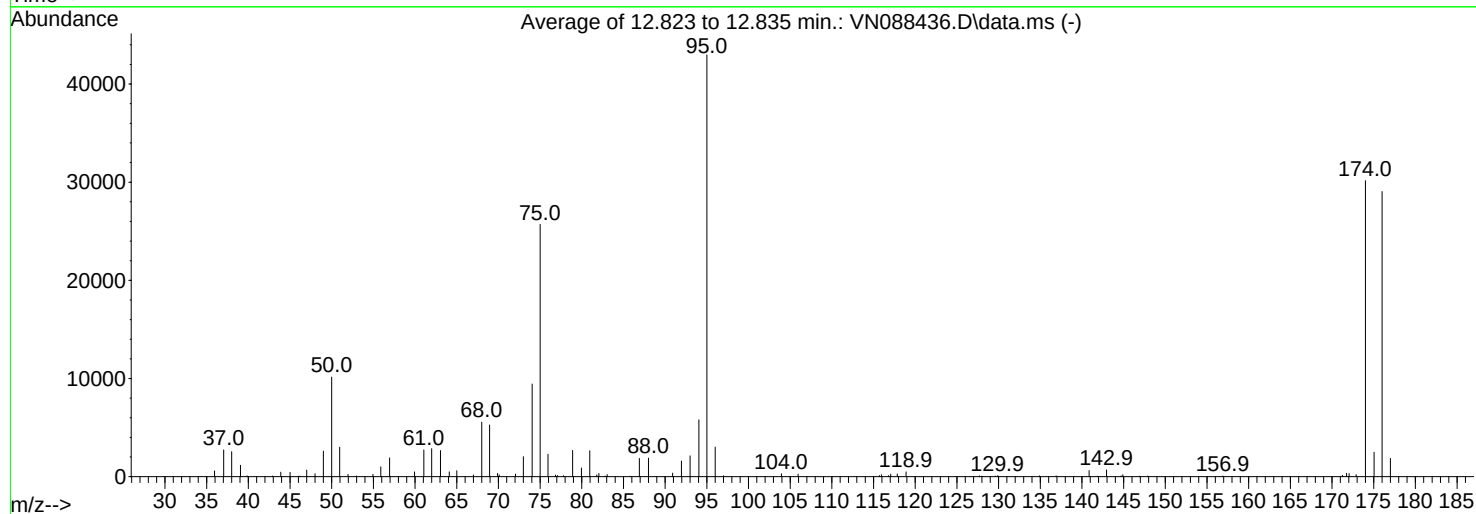
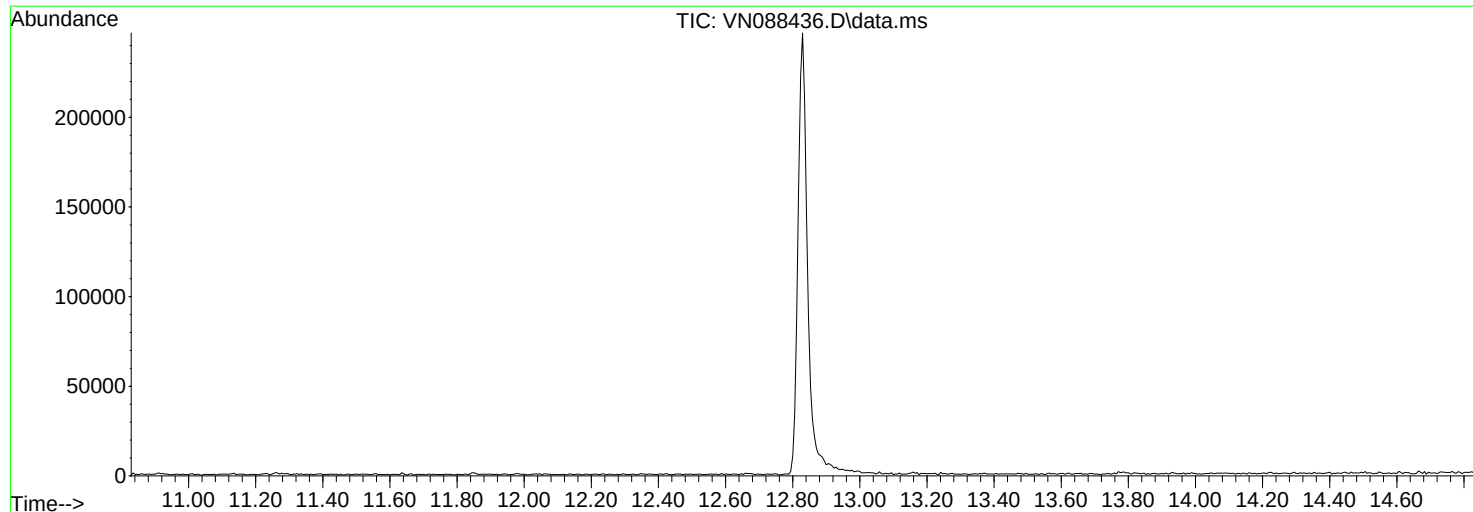
m/z	abund.	m/z	abund.	m/z	abund.	m/z	abund.
281.30	53						

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN120425\
Data File : VN088436.D
Acq On : 04 Dec 2025 11:03
Operator : JC\MD
Sample : BFB
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
BFB

Integration File: RTEINT.P

Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N112425W.M
Title : SW846 8260
Last Update : Tue Nov 25 01:34:52 2025



AutoFind: Scans 1848, 1849, 1850; Background Corrected with Scan 1839

Target	Rel. to	Lower	Upper	Rel.	Raw	Result
Mass	Mass	Limit%	Limit%	Abn%	Abn	Pass/Fail
50	95	15	40	23.6	10147	PASS
75	95	30	60	59.8	25707	PASS
95	95	100	100	100.0	42976	PASS
96	95	5	9	7.0	3020	PASS
173	174	0.00	2	0.7	221	PASS
174	95	50	100	70.2	30165	PASS
175	174	5	9	8.3	2492	PASS
176	174	95	101	96.3	29056	PASS
177	176	5	9	6.4	1867	PASS

BFB

Modified:subtracted

m/z	abund.	m/z	abund.	m/z	abund.	m/z	abund.
35.95	584	48.00	287	62.00	2852	74.05	943
37.05	2726	49.00	2602	63.05	2665	75.00	2570
38.00	2546	50.00	10147	64.10	494	75.95	228
39.05	1160	50.95	2998	65.00	605	76.85	17
39.85	81	51.95	228	67.00	181	77.10	110
42.90	65	53.00	65	68.00	5547	77.80	94
43.90	456	54.95	241	68.95	5266	78.90	2668
45.00	436	55.90	989	69.90	333	79.95	879
46.00	52	56.95	1910	70.10	186	80.95	2648
46.20	51	59.95	491	72.05	264	81.80	181
47.00	668	61.05	2720	73.00	2047	82.05	329

Average of 12.823 to 12.835 min.: VN088436.D\data.ms

BFB

Modified:subtracted

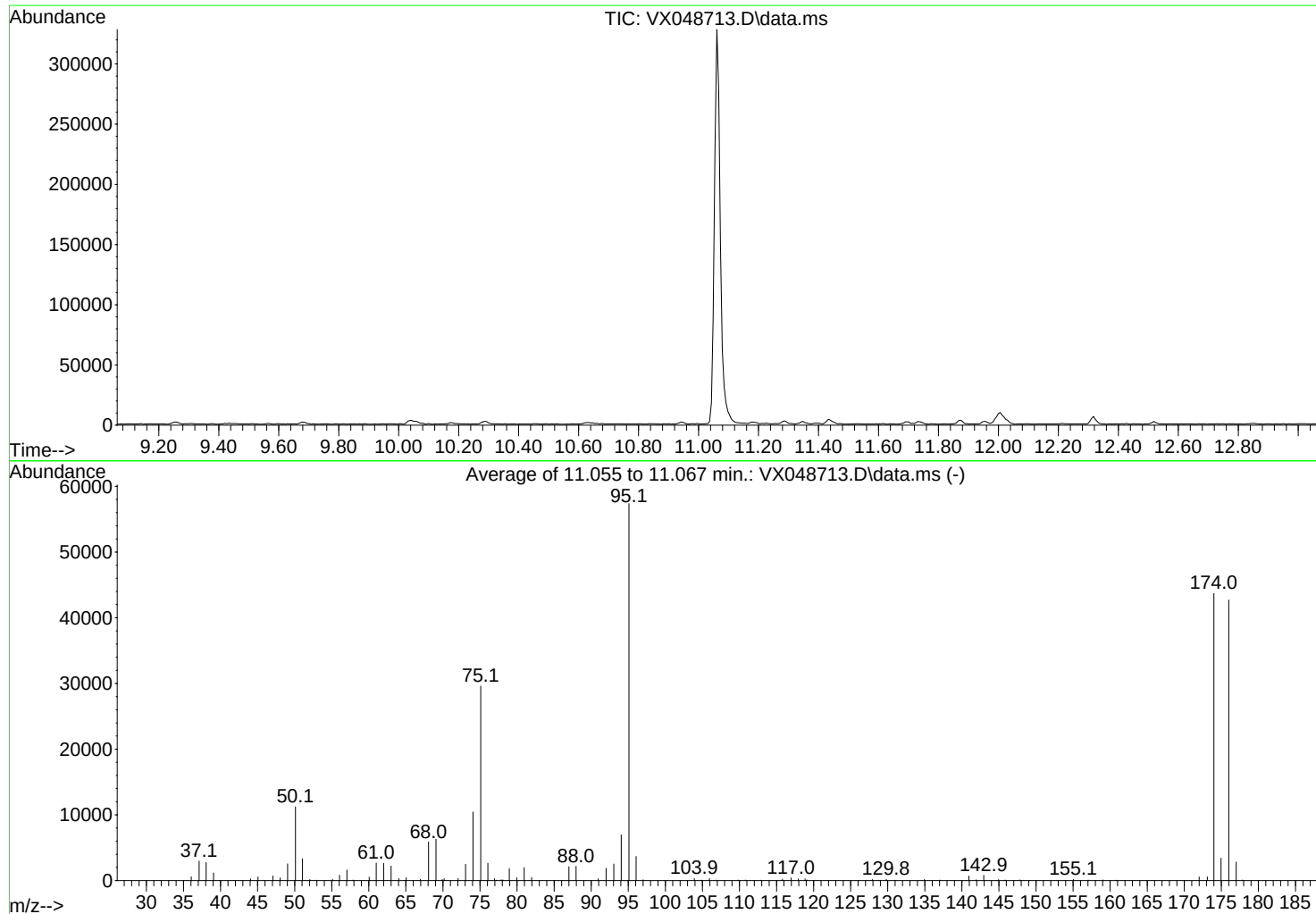
m/z	abund.	m/z	abund.	m/z	abund.	m/z	abund.
82.80	60	97.00	99	129.60	63	156.90	69
83.05	210	103.95	288	129.90	95	171.25	133
85.80	57	105.95	255	134.90	84	171.75	339
86.90	1861	115.70	111	136.90	80	172.05	316
88.00	1886	115.95	198	140.85	619	172.90	221
90.90	359	116.80	99	141.90	54	174.00	30165
91.95	1605	117.05	260	142.95	686	175.05	2492
93.00	2116	117.85	283	144.70	81	176.00	29056
94.05	5785	118.90	482	144.90	199	177.00	1867
95.00	42976	127.70	81	146.90	53		
96.00	3020	128.00	53	147.90	63		

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120425\
Data File : VX048713.D
Acq On : 04 Dec 2025 08:15
Operator : JC/MD
Sample : BFB
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
BFB

Integration File: RTEINT.P

Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X120425W.M
Title : SW846 8260
Last Update : Fri Dec 05 03:27:34 2025



AutoFind: Scans 1635, 1636, 1637; Background Corrected with Scan 1628

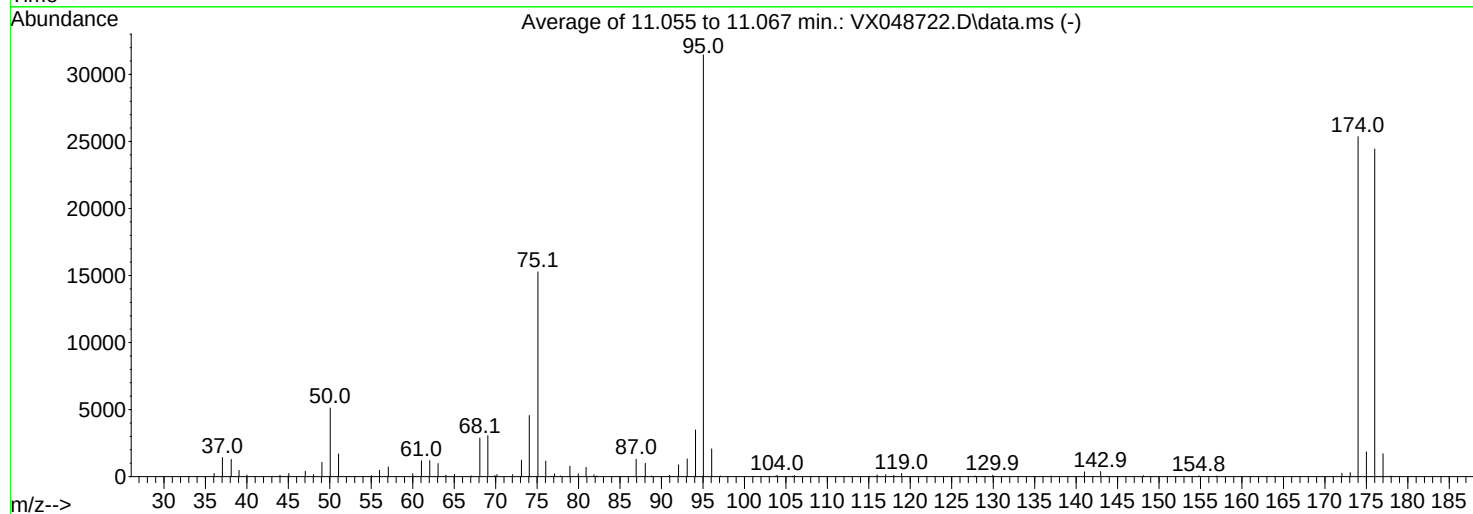
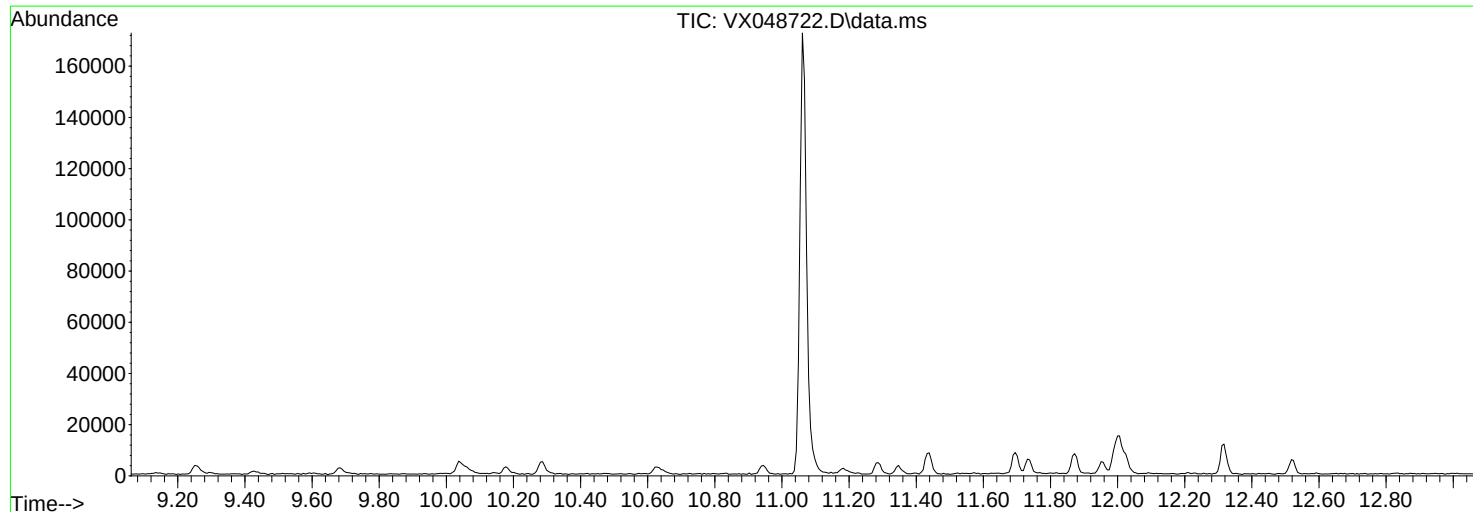
Target	Rel. to	Lower	Upper	Rel.	Raw	Result
Mass	Mass	Limit%	Limit%	Abn%	Abn	Pass/Fail
50	95	15	40	19.6	11214	PASS
75	95	30	60	51.6	29608	PASS
95	95	100	100	100.0	57360	PASS
96	95	5	9	6.4	3677	PASS
173	174	0.00	2	1.3	568	PASS
174	95	50	100	76.2	43696	PASS
175	174	5	9	7.8	3397	PASS
176	174	95	101	97.7	42696	PASS
177	176	5	9	6.6	2825	PASS

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120525\
Data File : VX048722.D
Acq On : 05 Dec 2025 08:14
Operator : JC/MD
Sample : BFB
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
BFB

Integration File: RTEINT.P

Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X120425W.M
Title : SW846 8260
Last Update : Fri Dec 05 03:27:34 2025



AutoFind: Scans 1635, 1636, 1637; Background Corrected with Scan 1629

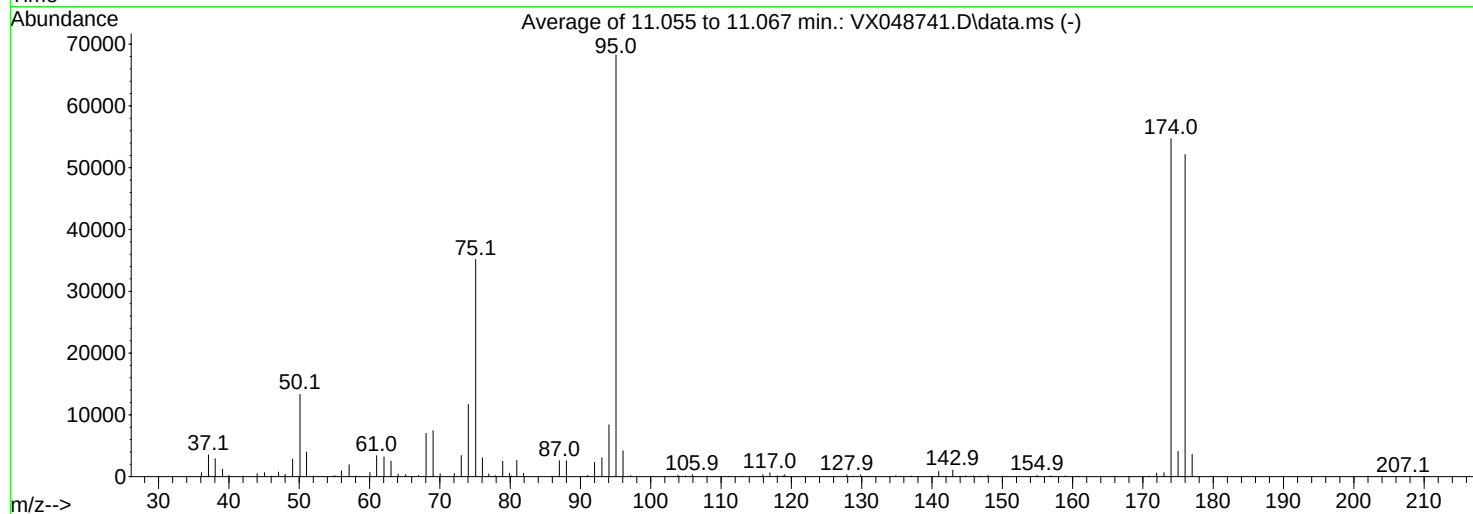
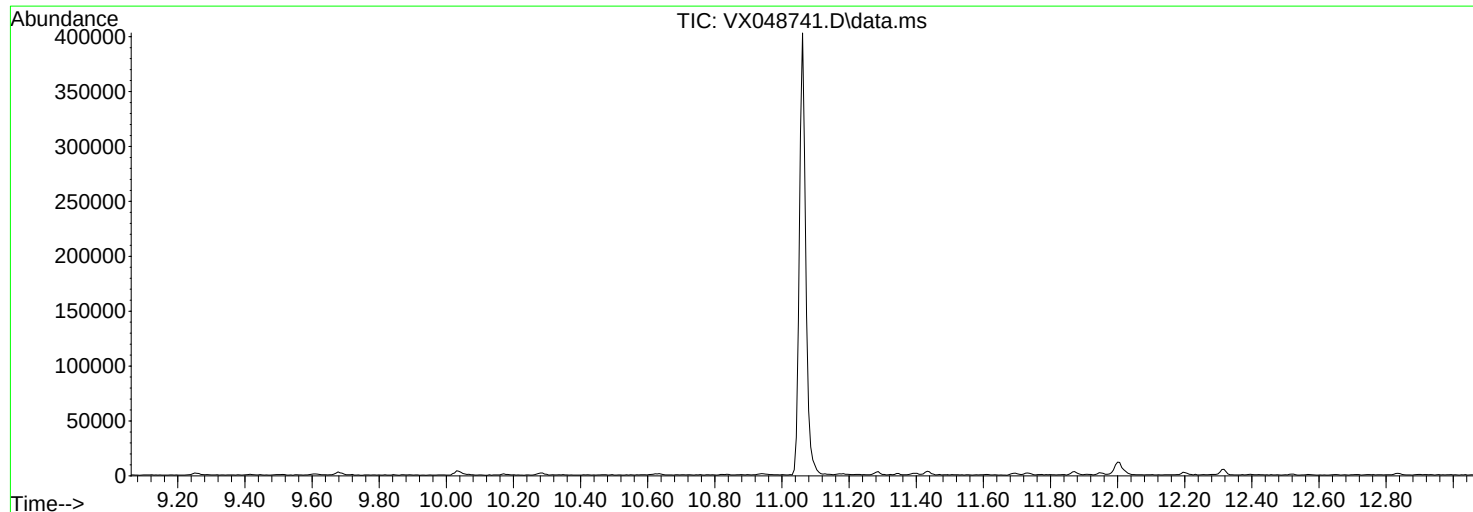
Target Mass	Rel. to Mass	Lower Limit%	Upper Limit%	Rel. Abn%	Raw Abn	Result Pass/Fail
50	95	15	40	16.3	5123	PASS
75	95	30	60	48.5	15274	PASS
95	95	100	100	100.0	31461	PASS
96	95	5	9	6.6	2062	PASS
173	174	0.00	2	1.2	294	PASS
174	95	50	100	80.6	25363	PASS
175	174	5	9	7.3	1854	PASS
176	174	95	101	96.4	24445	PASS
177	176	5	9	7.0	1707	PASS

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120825\
Data File : VX048741.D
Acq On : 08 Dec 2025 08:17
Operator : JC/MD
Sample : BFB
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
BFB

Integration File: RTEINT.P

Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X120425W.M
Title : SW846 8260
Last Update : Fri Dec 05 03:27:34 2025



AutoFind: Scans 1635, 1636, 1637; Background Corrected with Scan 1629

Target	Rel. to	Lower	Upper	Rel.	Raw	Result
Mass	Mass	Limit%	Limit%	Abn%	Abn	Pass/Fail
50	95	15	40	19.5	13323	PASS
75	95	30	60	51.5	35152	PASS
95	95	100	100	100.0	68272	PASS
96	95	5	9	6.2	4205	PASS
173	174	0.00	2	1.2	656	PASS
174	95	50	100	80.1	54717	PASS
175	174	5	9	7.5	4103	PASS
176	174	95	101	95.3	52133	PASS
177	176	5	9	7.0	3634	PASS



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

Report of Analysis

Client: JACOBS Engineering Group, Inc.
Project: Former Schlumberger STC PTC Site D3868221
Client Sample ID: VN1204WBL01
Lab Sample ID: VN1204WBL01
Analytical Method: 8260D
Sample Wt/Vol: 5 mL

Level: LOW
Final Vol: 5000 uL

Date Collected:
Date Received:
SDG No.: Q3757
Matrix: Water
% Solid: 0
Test: VOCMS Group3

CAS Number	Parameter	Conc.	Qua.	DF	MDL	LOQ / CRQL	Units	Date Ana.	BatchID
TARGETS									
75-01-4	Vinyl Chloride	0.26	U	1	0.26	1.00	ug/L	12/04/25 12:38	VN120425
75-35-4	1,1-Dichloroethene	0.23	U	1	0.23	1.00	ug/L	12/04/25 12:38	VN120425
75-34-3	1,1-Dichloroethane	0.23	U	1	0.23	1.00	ug/L	12/04/25 12:38	VN120425
156-59-2	cis-1,2-Dichloroethene	0.19	U	1	0.19	1.00	ug/L	12/04/25 12:38	VN120425
71-55-6	1,1,1-Trichloroethane	0.20	U	1	0.20	1.00	ug/L	12/04/25 12:38	VN120425
71-43-2	Benzene	0.15	U	1	0.15	1.00	ug/L	12/04/25 12:38	VN120425
107-06-2	1,2-Dichloroethane	0.22	U	1	0.22	1.00	ug/L	12/04/25 12:38	VN120425
79-01-6	Trichloroethene	0.090	U	1	0.090	1.00	ug/L	12/04/25 12:38	VN120425
79-00-5	1,1,2-Trichloroethane	0.21	U	1	0.21	1.00	ug/L	12/04/25 12:38	VN120425
127-18-4	Tetrachloroethene	0.23	U	1	0.23	1.00	ug/L	12/04/25 12:38	VN120425
SURROGATES									
17060-07-0	1,2-Dichloroethane-d4	61.7			70 (74) - 130 (125)	123%	SPK: 50		
1868-53-7	Dibromofluoromethane	50.8			70 (75) - 130 (124)	102%	SPK: 50		
2037-26-5	Toluene-d8	46.7			70 (86) - 130 (113)	93%	SPK: 50		
460-00-4	4-Bromofluorobenzene	49.9			70 (77) - 130 (121)	100%	SPK: 50		
INTERNAL STANDARDS									
		Area Count							
363-72-4	Pentafluorobenzene	237000							
540-36-3	1,4-Difluorobenzene	557000							
3114-55-4	Chlorobenzene-d5	556000							
3855-82-1	1,4-Dichlorobenzene-d4	258000							

U = Not Detected

LOQ = Limit of Quantitation

MDL = Method Detection Limit

LOD = Limit of Detection

E = Value Exceeds Calibration Range

Q = indicates LCS control criteria did not meet requirements

M = MS/MSD acceptance criteria did not meet requirements

J = Estimated Value

B = Analyte Found in Associated Method Blank

N = Presumptive Evidence of a Compound

* = Values outside of QC limits

D = Dilution

() = Laboratory InHouse Limit

A = Aldol-Condensation Reaction Products

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN120425\
 Data File : VN088439.D
 Acq On : 04 Dec 2025 12:38
 Operator : JC\MD
 Sample : VN1204WBL01
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 VN1204WBL01

Quant Time: Dec 05 00:12:44 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N112425W.M
 Quant Title : SW846 8260
 QLast Update : Tue Nov 25 01:34:52 2025
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.206	168	236600	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.082	114	556948	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.841	117	555675	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.770	152	258091	50.000	ug/l	0.00

System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.559	65	275126	61.728	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	123.460%
35) Dibromofluoromethane	8.147	113	196090	50.805	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	101.620%
50) Toluene-d8	10.541	98	670134	46.664	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	93.320%
62) 4-Bromofluorobenzene	12.829	95	246023	49.929	ug/l	0.00
Spiked Amount	50.000	Range	77 - 121	Recovery	=	99.860%

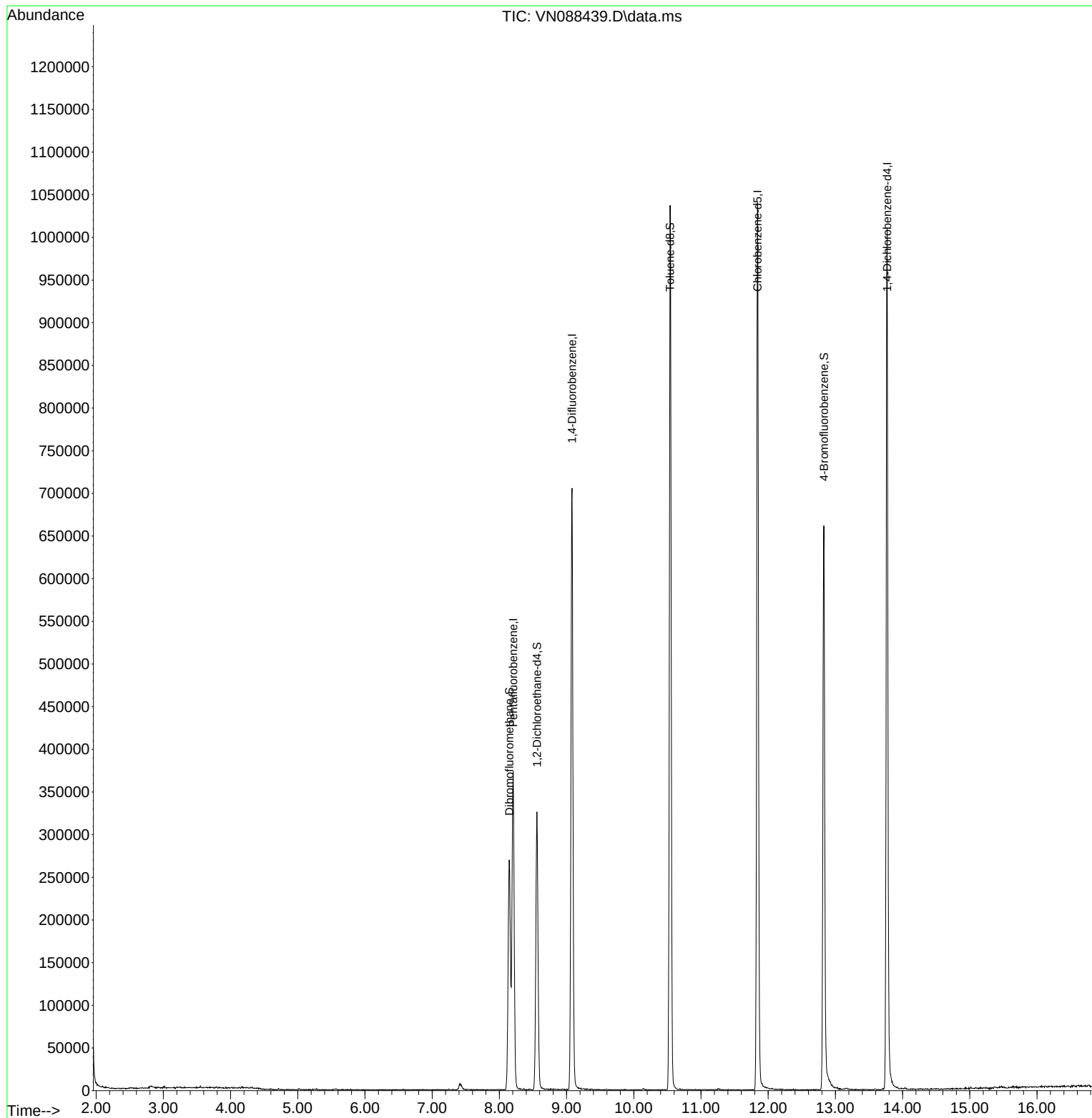
Target Compounds	Qvalue

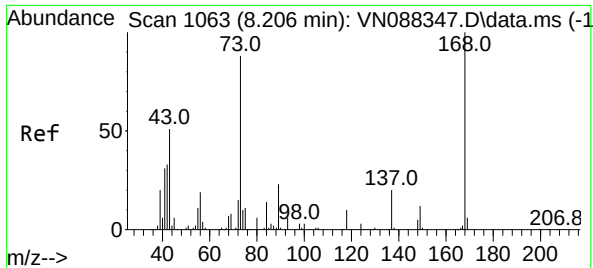
(#) = qualifier out of range (m) = manual integration (+) = signals summed

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN120425\
Data File : VN088439.D
Acq On : 04 Dec 2025 12:38
Operator : JC\MD
Sample : VN1204WBL01
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 5 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
VN1204WBL01

Quant Time: Dec 05 00:12:44 2025
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N112425W.M
Quant Title : SW846 8260
QLast Update : Tue Nov 25 01:34:52 2025
Response via : Initial Calibration

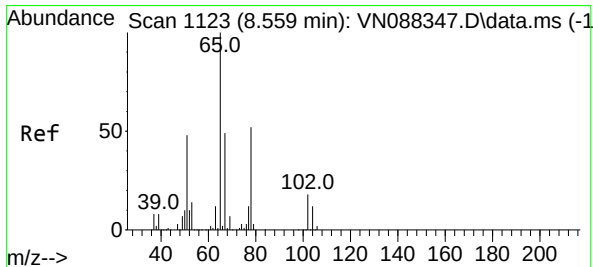
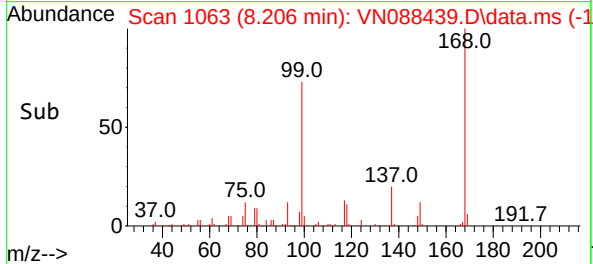
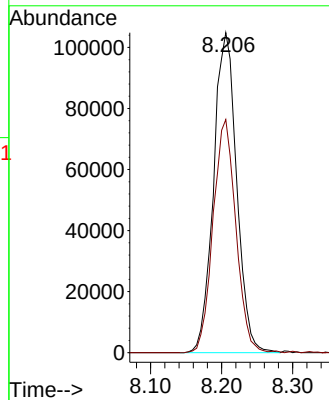
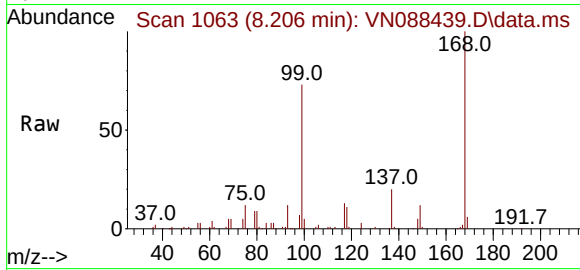




#1
Pentafluorobenzene
Concen: 50.000 ug/l
RT: 8.206 min Scan# 1063
Delta R.T. -0.000 min
Lab File: VN088439.D
Acq: 04 Dec 2025 12:38

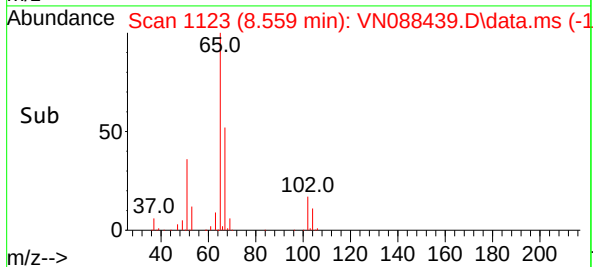
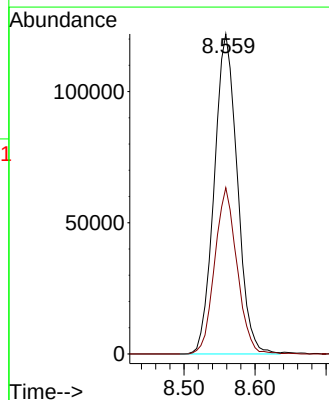
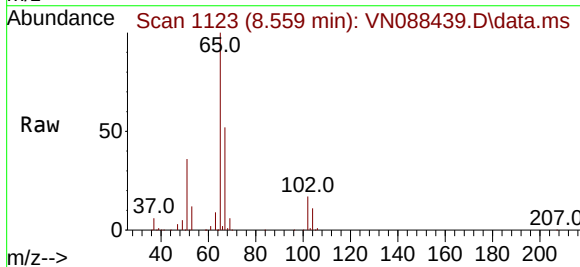
Instrument :
MSVOA_N
ClientSampleId :
VN1204WBL01

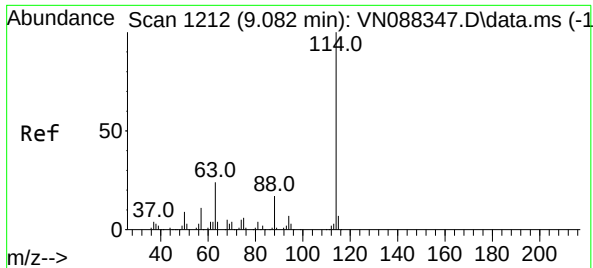
Tgt Ion:168 Resp: 236600
Ion Ratio Lower Upper
168 100
99 72.9 57.1 85.7



#33
1,2-Dichloroethane-d4
Concen: 61.728 ug/l
RT: 8.559 min Scan# 1123
Delta R.T. -0.000 min
Lab File: VN088439.D
Acq: 04 Dec 2025 12:38

Tgt Ion: 65 Resp: 275126
Ion Ratio Lower Upper
65 100
67 50.2 0.0 100.8





#34

1,4-Difluorobenzene

Concen: 50.000 ug/l

RT: 9.082 min Scan# 11

Delta R.T. -0.000 min

Lab File: VN088439.D

Acq: 04 Dec 2025 12:38

Instrument :

MSVOA_N

ClientSampleId :

VN1204WBL01

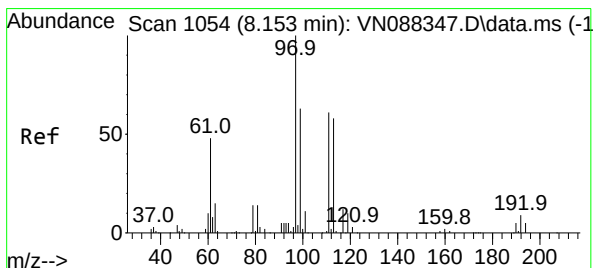
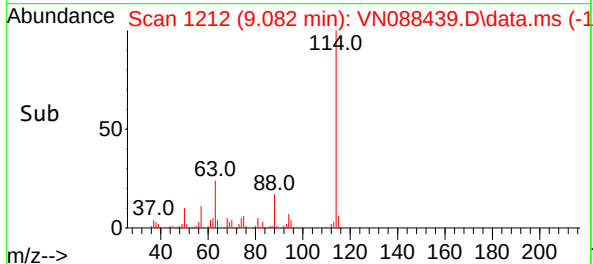
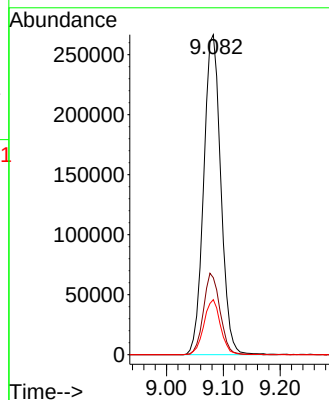
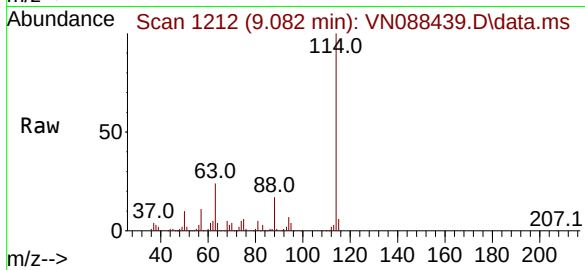
Tgt Ion:114 Resp: 556948

Ion Ratio Lower Upper

114 100

63 24.1 0.0 49.0

88 17.2 0.0 34.0



#35

Dibromofluoromethane

Concen: 50.805 ug/l

RT: 8.147 min Scan# 1053

Delta R.T. -0.006 min

Lab File: VN088439.D

Acq: 04 Dec 2025 12:38

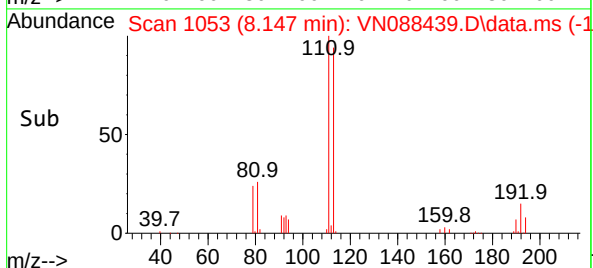
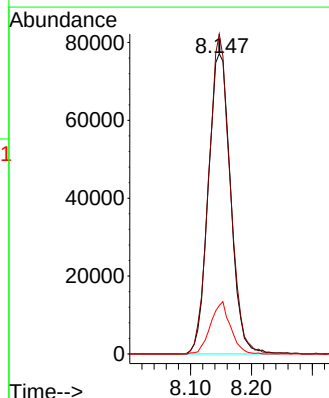
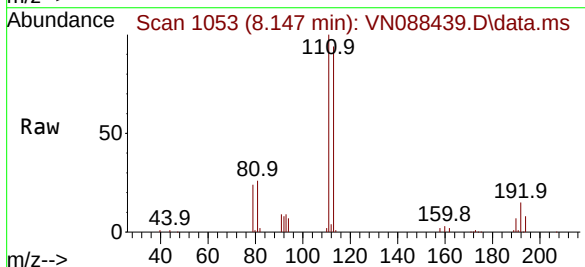
Tgt Ion:113 Resp: 196090

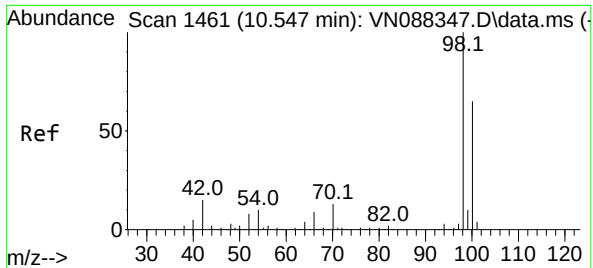
Ion Ratio Lower Upper

113 100

111 102.2 82.5 123.7

192 15.6 12.8 19.2

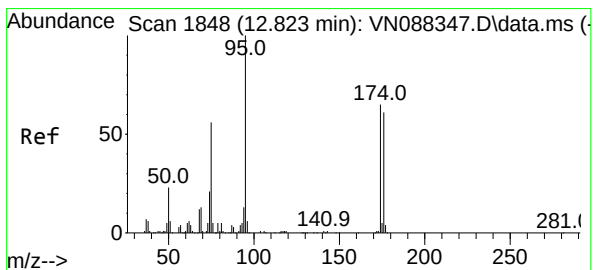
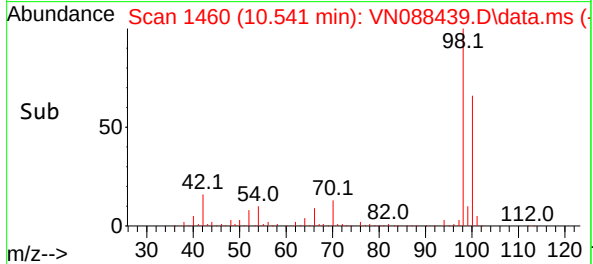
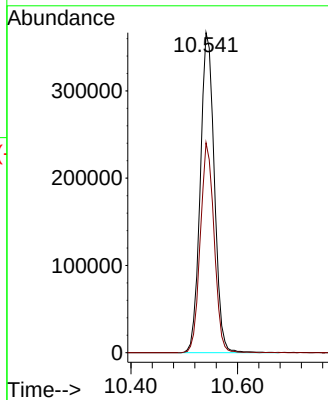
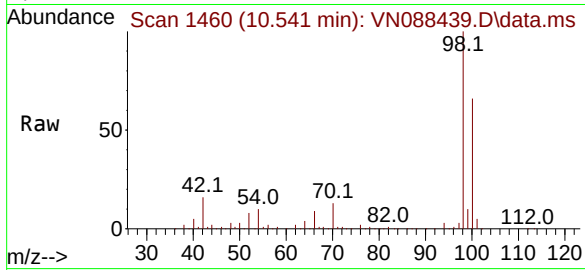




#50
Toluene-d8
Concen: 46.664 ug/l
RT: 10.541 min Scan# 1460
Delta R.T. -0.006 min
Lab File: VN088439.D
Acq: 04 Dec 2025 12:38

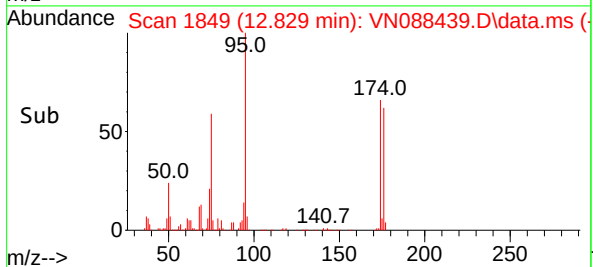
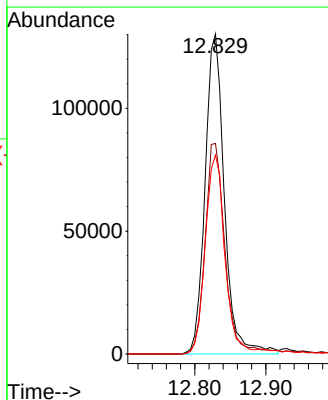
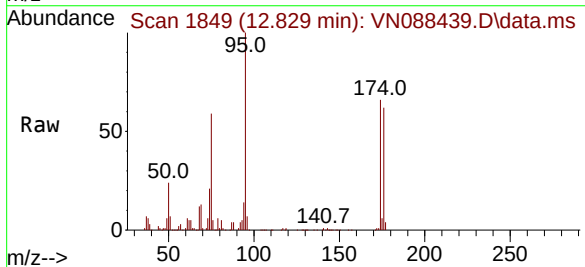
Instrument :
MSVOA_N
ClientSampleId :
VN1204WBL01

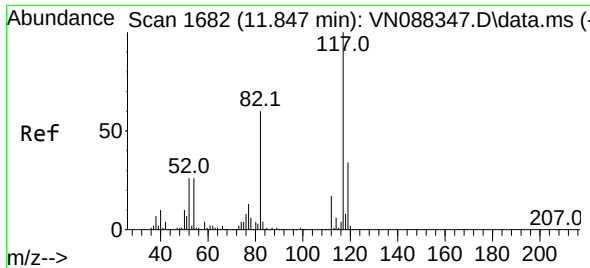
Tgt Ion: 98 Resp: 670134
Ion Ratio Lower Upper
98 100
100 64.4 52.2 78.2



#62
4-Bromofluorobenzene
Concen: 49.929 ug/l
RT: 12.829 min Scan# 1849
Delta R.T. 0.006 min
Lab File: VN088439.D
Acq: 04 Dec 2025 12:38

Tgt Ion: 95 Resp: 246023
Ion Ratio Lower Upper
95 100
174 68.9 0.0 139.0
176 62.6 0.0 132.4

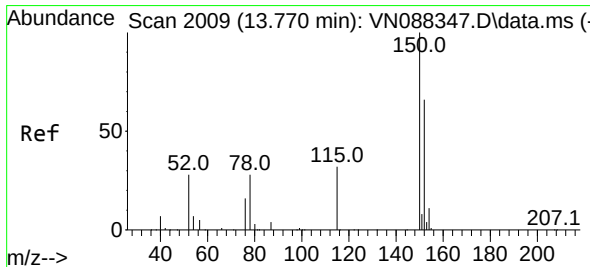
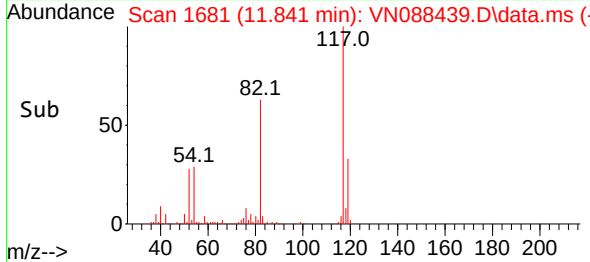
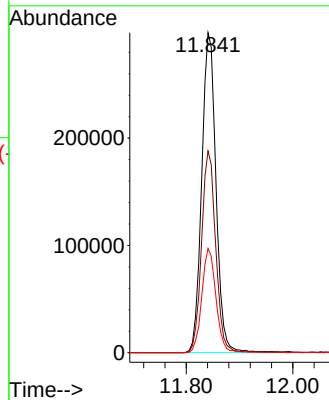
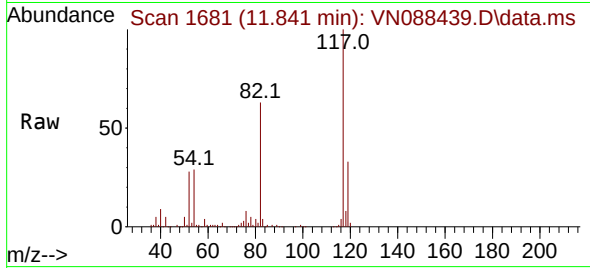




#63
Chlorobenzene-d5
Concen: 50.000 ug/l
RT: 11.841 min Scan# 117
Delta R.T. -0.006 min
Lab File: VN088439.D
Acq: 04 Dec 2025 12:38

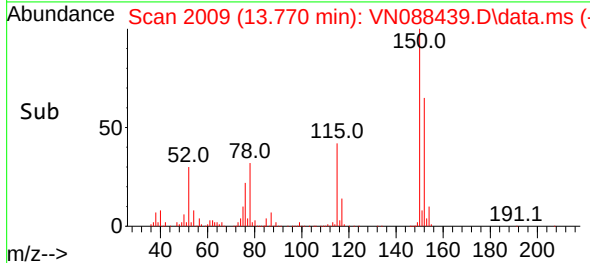
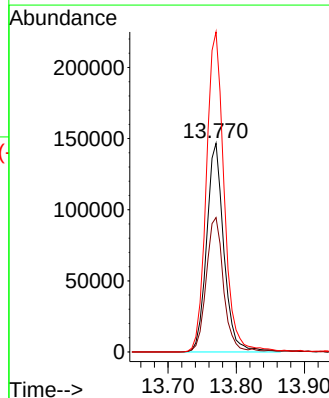
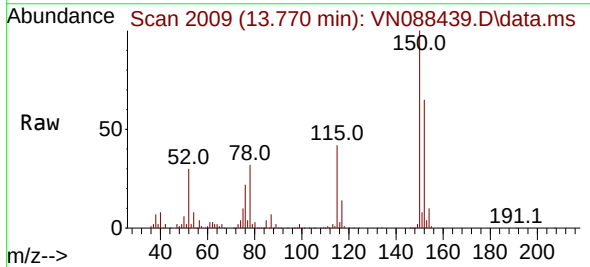
Instrument :
MSVOA_N
ClientSampleId :
VN1204WBL01

Tgt Ion	Ratio	Lower	Upper
117	100		
82	63.0	47.8	71.8
119	32.6	26.9	40.3



#72
1,4-Dichlorobenzene-d4
Concen: 50.000 ug/l
RT: 13.770 min Scan# 2009
Delta R.T. -0.000 min
Lab File: VN088439.D
Acq: 04 Dec 2025 12:38

Tgt Ion	Ratio	Lower	Upper
152	100		
115	64.9	33.0	98.9
150	158.8	0.0	349.0





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

Report of Analysis

Client: JACOBS Engineering Group, Inc.
Project: Former Schlumberger STC PTC Site D3868221
Client Sample ID: VX1205WBL01
Lab Sample ID: VX1205WBL01
Analytical Method: 8260D
Sample Wt/Vol: 5 mL

Level: LOW
Final Vol: 5000 uL

Date Collected:
Date Received:
SDG No.: Q3757
Matrix: Water
% Solid: 0
Test: VOCMS Group3

CAS Number	Parameter	Conc.	Qua.	DF	MDL	LOQ / CRQL	Units	Date Ana.	BatchID
TARGETS									
75-01-4	Vinyl Chloride	0.26	U	1	0.26	1.00	ug/L	12/05/25 10:34	VX120525
75-35-4	1,1-Dichloroethene	0.23	U	1	0.23	1.00	ug/L	12/05/25 10:34	VX120525
75-34-3	1,1-Dichloroethane	0.23	U	1	0.23	1.00	ug/L	12/05/25 10:34	VX120525
156-59-2	cis-1,2-Dichloroethene	0.19	U	1	0.19	1.00	ug/L	12/05/25 10:34	VX120525
71-55-6	1,1,1-Trichloroethane	0.20	U	1	0.20	1.00	ug/L	12/05/25 10:34	VX120525
71-43-2	Benzene	0.15	U	1	0.15	1.00	ug/L	12/05/25 10:34	VX120525
107-06-2	1,2-Dichloroethane	0.22	U	1	0.22	1.00	ug/L	12/05/25 10:34	VX120525
79-01-6	Trichloroethene	0.090	U	1	0.090	1.00	ug/L	12/05/25 10:34	VX120525
79-00-5	1,1,2-Trichloroethane	0.21	U	1	0.21	1.00	ug/L	12/05/25 10:34	VX120525
127-18-4	Tetrachloroethene	0.23	U	1	0.23	1.00	ug/L	12/05/25 10:34	VX120525
SURROGATES									
17060-07-0	1,2-Dichloroethane-d4	50.6			70 (74) - 130 (125)	101%	SPK: 50		
1868-53-7	Dibromofluoromethane	49.7			70 (75) - 130 (124)	99%	SPK: 50		
2037-26-5	Toluene-d8	50.8			70 (86) - 130 (113)	102%	SPK: 50		
460-00-4	4-Bromofluorobenzene	60.5			70 (77) - 130 (121)	121%	SPK: 50		
INTERNAL STANDARDS									
		Area Count							
363-72-4	Pentafluorobenzene	193000							
540-36-3	1,4-Difluorobenzene	291000							
3114-55-4	Chlorobenzene-d5	294000							
3855-82-1	1,4-Dichlorobenzene-d4	131000							

U = Not Detected

LOQ = Limit of Quantitation

MDL = Method Detection Limit

LOD = Limit of Detection

E = Value Exceeds Calibration Range

Q = indicates LCS control criteria did not meet requirements

M = MS/MSD acceptance criteria did not meet requirements

J = Estimated Value

B = Analyte Found in Associated Method Blank

N = Presumptive Evidence of a Compound

* = Values outside of QC limits

D = Dilution

() = Laboratory InHouse Limit

A = Aldol-Condensation Reaction Products

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120525\
 Data File : VX048725.D
 Acq On : 05 Dec 2025 10:34
 Operator : JC/MD
 Sample : VX1205WBL01
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 VX1205WBL01

Quant Time: Dec 05 21:59:18 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X120425W.M
 Quant Title : SW846 8260
 QLast Update : Fri Dec 05 03:27:34 2025
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.532	168	193187	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.739	114	291151	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.037	117	294164	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.006	152	130501	50.000	ug/l	0.00

System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.928	65	131204	50.601	ug/l	-0.01
Spiked Amount	50.000	Range	78 - 117	Recovery	=	101.200%
35) Dibromofluoromethane	5.367	113	99604	49.686	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	99.380%
50) Toluene-d8	8.629	98	356802	50.775	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	101.560%
62) 4-Bromofluorobenzene	11.061	95	146671	60.529	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	121.060%

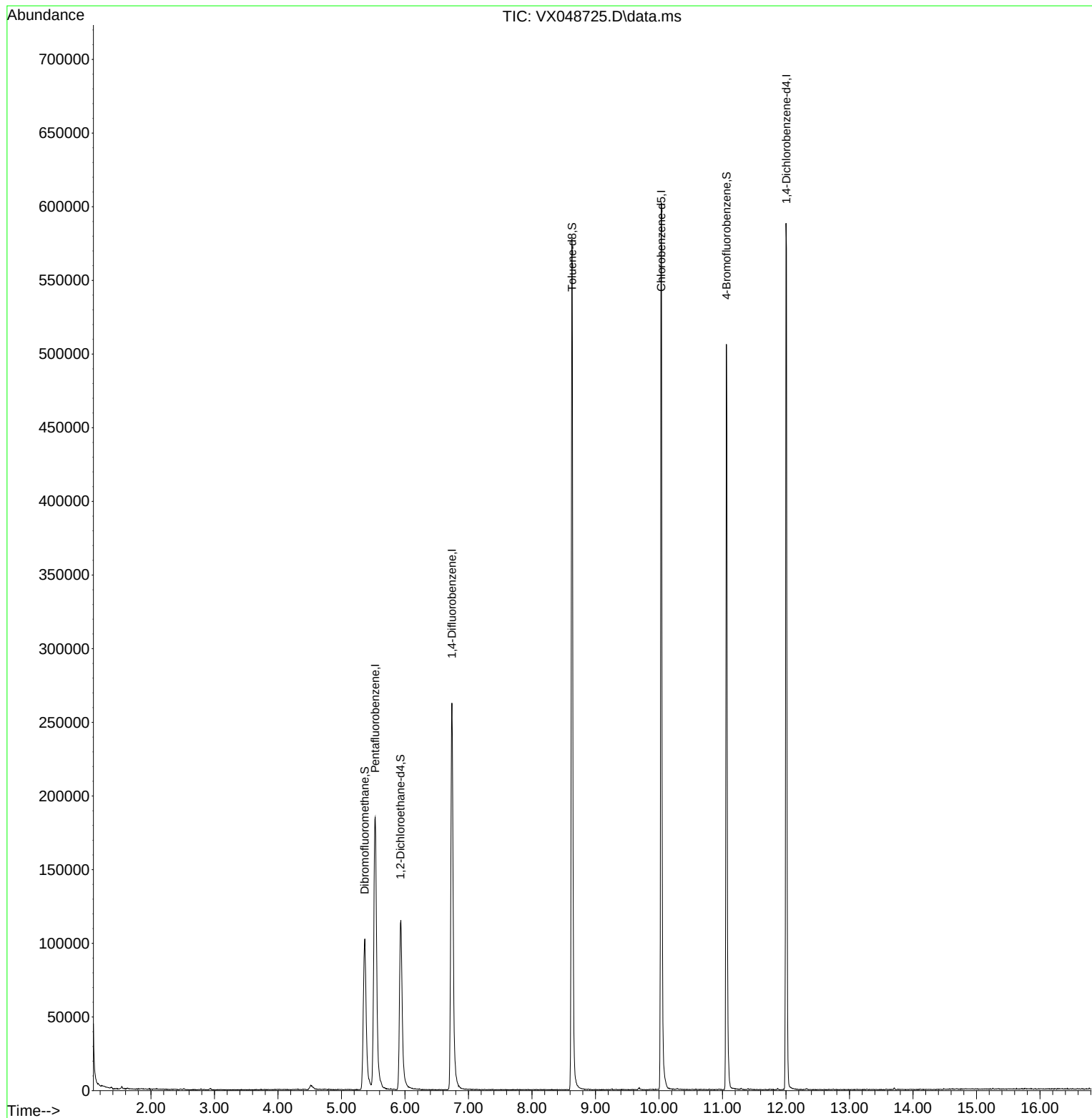
Target Compounds	Qvalue

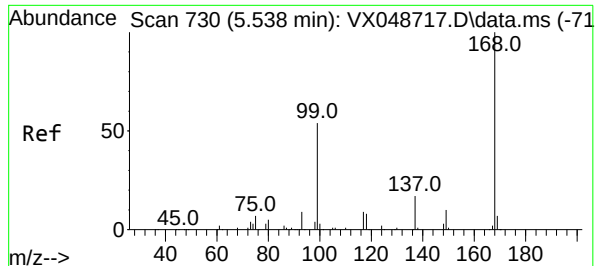
(#) = qualifier out of range (m) = manual integration (+) = signals summed

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120525\
Data File : VX048725.D
Acq On : 05 Dec 2025 10:34
Operator : JC/MD
Sample : VX1205WBL01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 4 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
VX1205WBL01

Quant Time: Dec 05 21:59:18 2025
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X120425W.M
Quant Title : SW846 8260
QLast Update : Fri Dec 05 03:27:34 2025
Response via : Initial Calibration

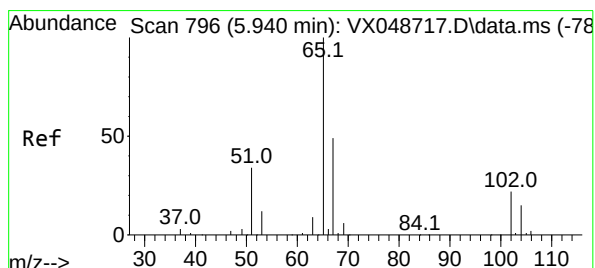
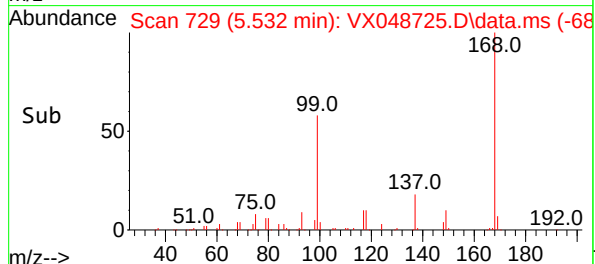
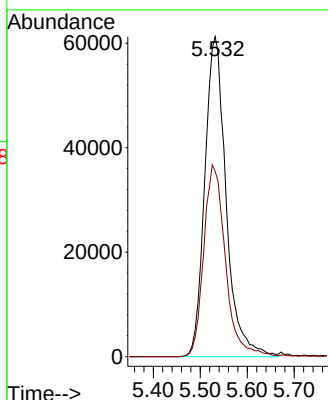
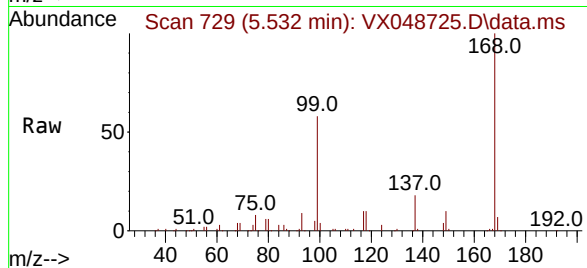




#1
Pentafluorobenzene
Concen: 50.000 ug/l
RT: 5.532 min Scan# 71
Delta R.T. -0.006 min
Lab File: VX048725.D
Acq: 05 Dec 2025 10:34

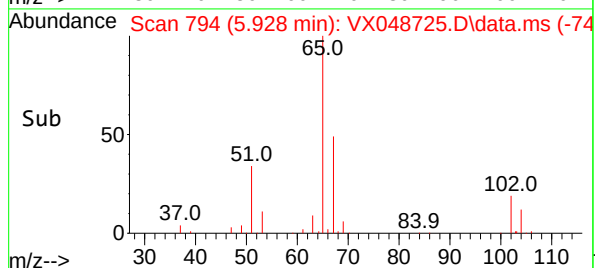
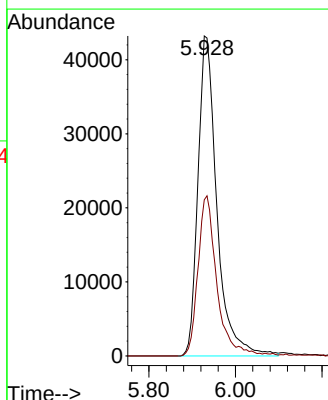
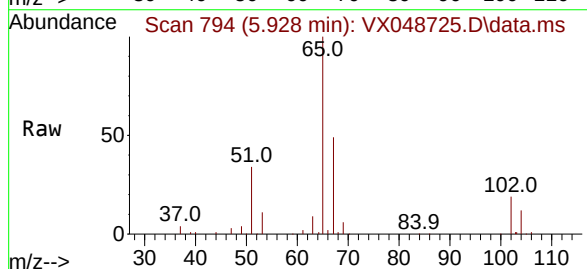
Instrument :
MSVOA_X
ClientSampleId :
VX1205WBL01

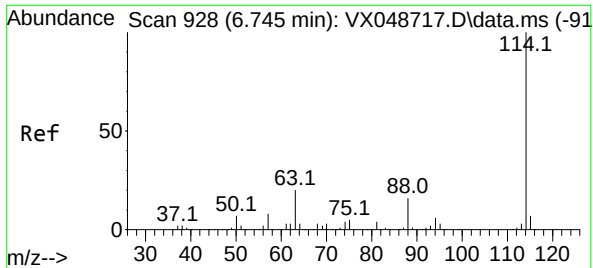
Tgt Ion:168 Resp: 193187
Ion Ratio Lower Upper
168 100
99 57.9 44.2 66.4



#33
1,2-Dichloroethane-d4
Concen: 50.601 ug/l
RT: 5.928 min Scan# 794
Delta R.T. -0.012 min
Lab File: VX048725.D
Acq: 05 Dec 2025 10:34

Tgt Ion: 65 Resp: 131204
Ion Ratio Lower Upper
65 100
67 49.6 0.0 107.4





#34

1,4-Difluorobenzene

Concen: 50.000 ug/l

RT: 6.739 min Scan# 91

Delta R.T. -0.006 min

Lab File: VX048725.D

Acq: 05 Dec 2025 10:34

Instrument :

MSVOA_X

ClientSampleId :

VX1205WBL01

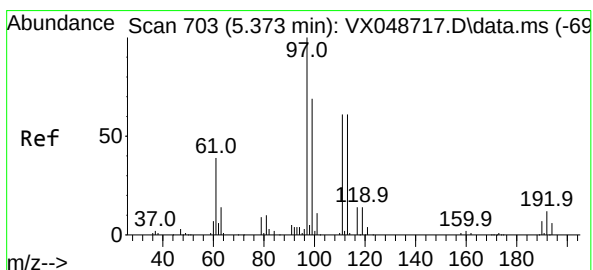
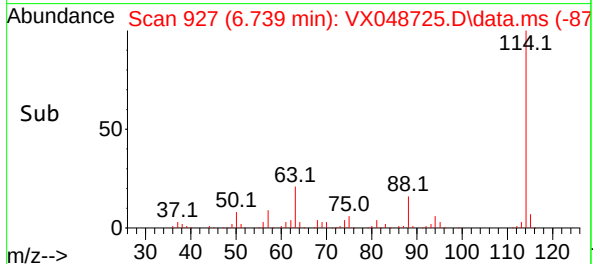
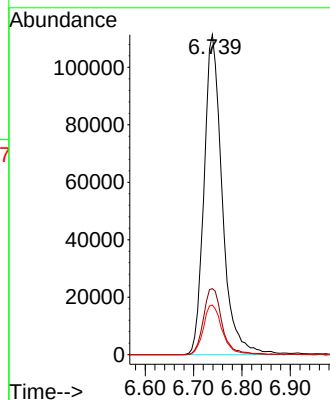
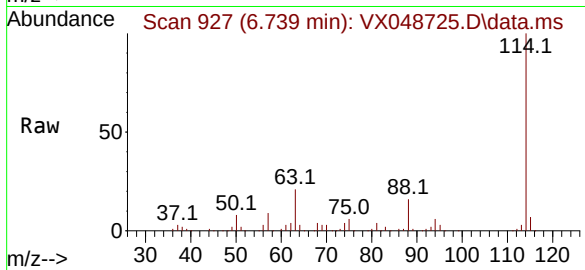
Tgt Ion:114 Resp: 291151

Ion Ratio Lower Upper

114 100

63 20.6 0.0 39.0

88 15.5 0.0 31.8



#35

Dibromofluoromethane

Concen: 49.686 ug/l

RT: 5.367 min Scan# 702

Delta R.T. -0.006 min

Lab File: VX048725.D

Acq: 05 Dec 2025 10:34

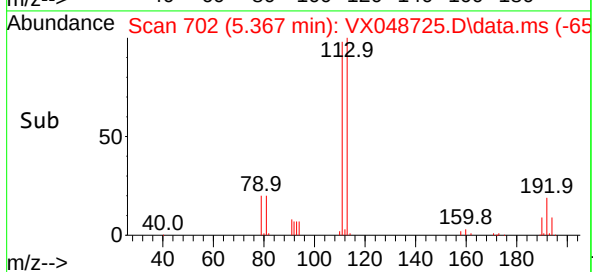
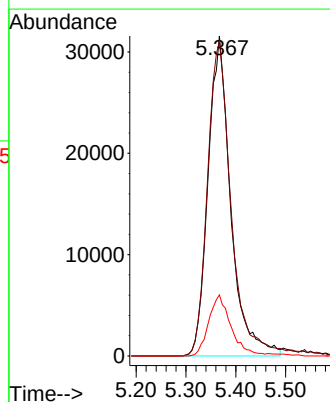
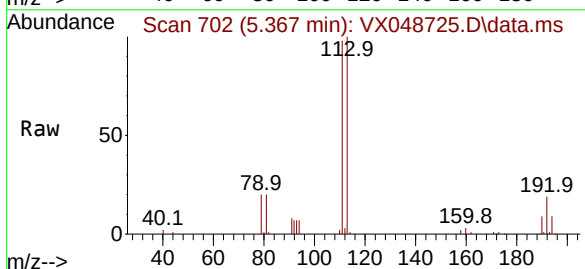
Tgt Ion:113 Resp: 99604

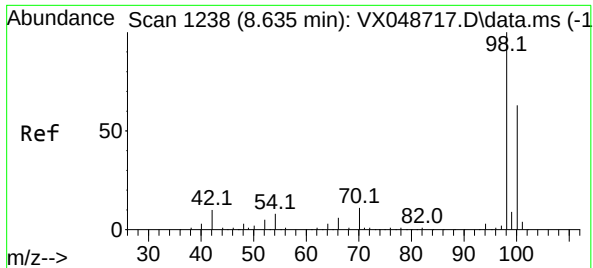
Ion Ratio Lower Upper

113 100

111 103.4 79.5 119.3

192 18.7 16.1 24.1

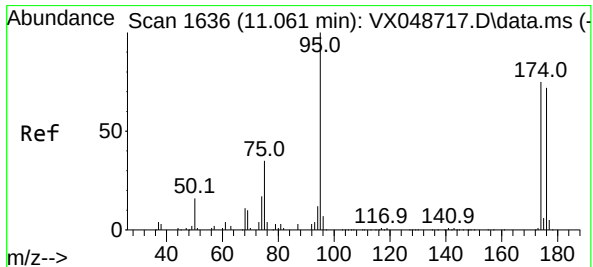
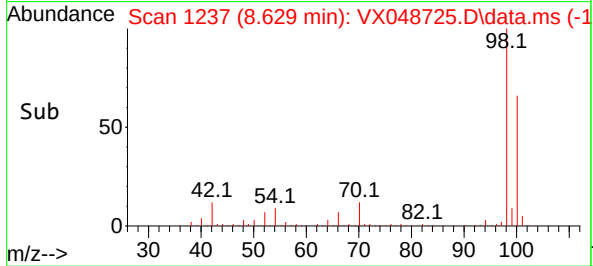
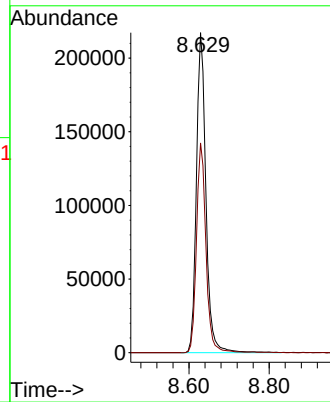
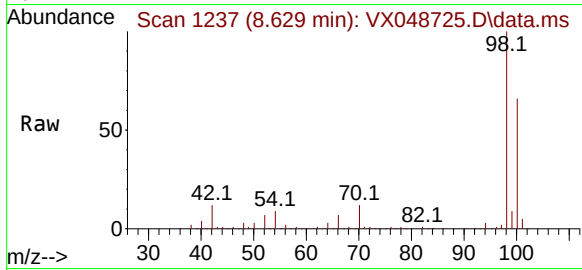




#50
Toluene-d8
Concen: 50.775 ug/l
RT: 8.629 min Scan# 1238
Delta R.T. -0.006 min
Lab File: VX048725.D
Acq: 05 Dec 2025 10:34

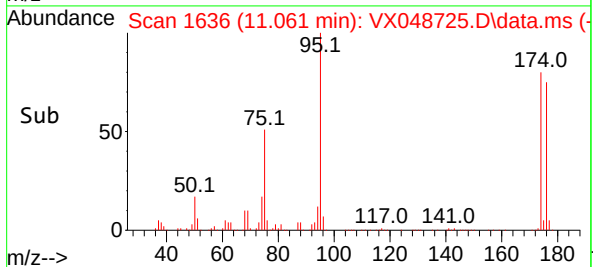
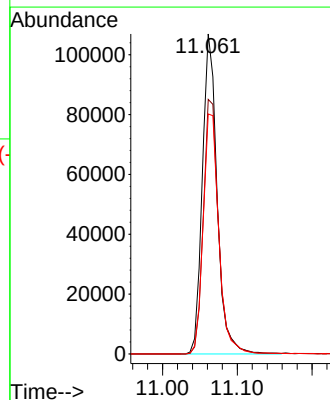
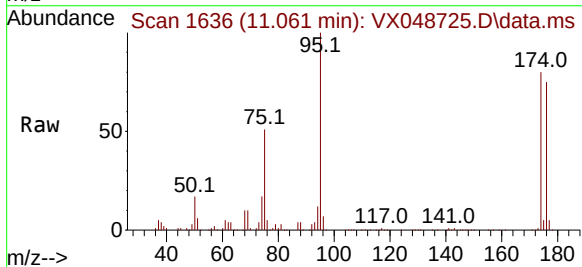
Instrument :
MSVOA_X
ClientSampleId :
VX1205WBL01

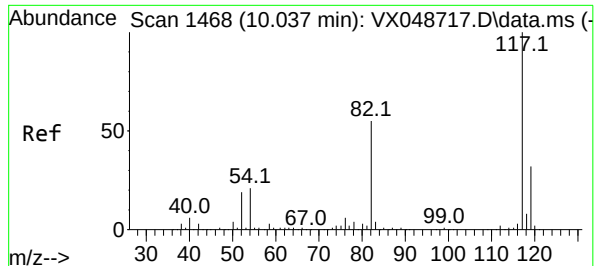
Tgt Ion: 98 Resp: 356802
Ion Ratio Lower Upper
98 100
100 65.3 53.4 80.0



#62
4-Bromofluorobenzene
Concen: 60.529 ug/l
RT: 11.061 min Scan# 1636
Delta R.T. 0.000 min
Lab File: VX048725.D
Acq: 05 Dec 2025 10:34

Tgt Ion: 95 Resp: 146671
Ion Ratio Lower Upper
95 100
174 82.3 0.0 157.8
176 79.8 0.0 154.0





#63

Chlorobenzene-d5

Concen: 50.000 ug/l

RT: 10.037 min Scan# 1468

Delta R.T. 0.000 min

Lab File: VX048725.D

Acq: 05 Dec 2025 10:34

Instrument :

MSVOA_X

ClientSampleId :

VX1205WBL01

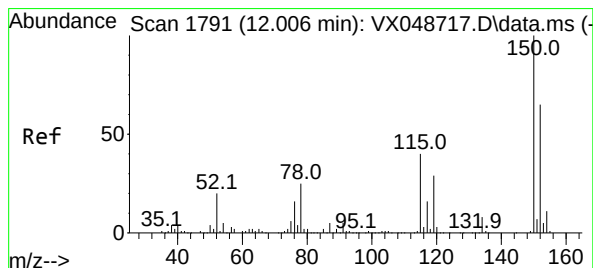
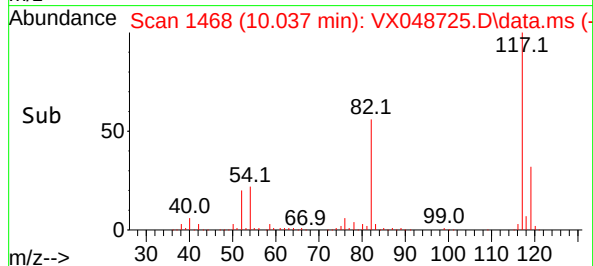
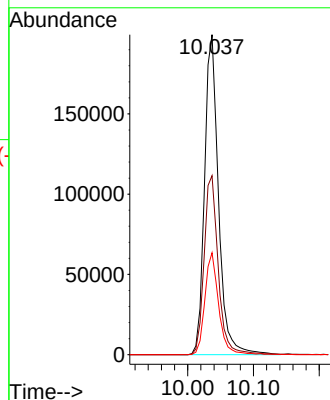
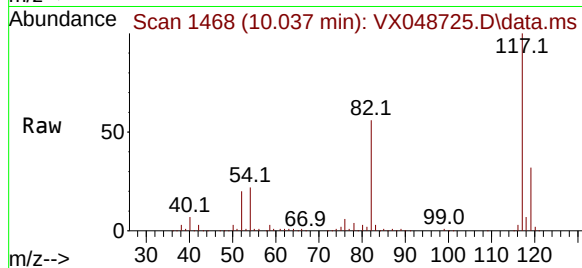
Tgt Ion:117 Resp: 294164

Ion Ratio Lower Upper

117 100

82 56.0 44.1 66.1

119 31.9 25.2 37.8



#72

1,4-Dichlorobenzene-d4

Concen: 50.000 ug/l

RT: 12.006 min Scan# 1791

Delta R.T. 0.000 min

Lab File: VX048725.D

Acq: 05 Dec 2025 10:34

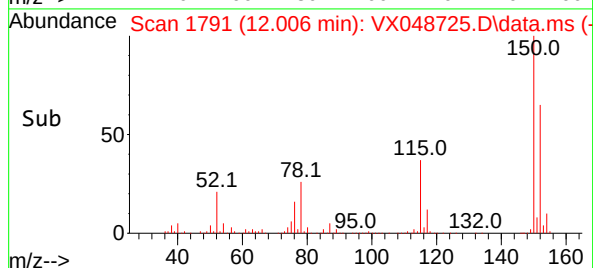
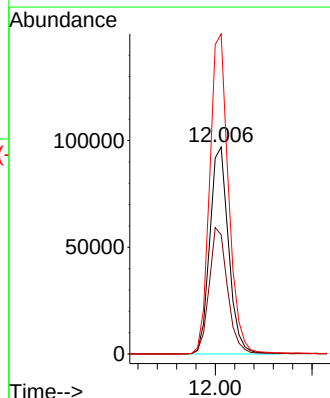
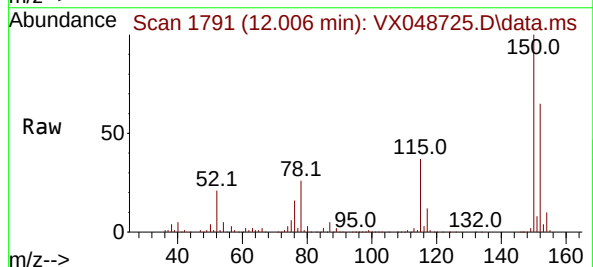
Tgt Ion:152 Resp: 130501

Ion Ratio Lower Upper

152 100

115 60.2 42.1 126.4

150 155.0 0.0 347.8





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

Report of Analysis

Client: JACOBS Engineering Group, Inc.
Project: Former Schlumberger STC PTC Site D3868221
Client Sample ID: VX1208WBL01
Lab Sample ID: VX1208WBL01
Analytical Method: 8260D
Sample Wt/Vol: 5 mL

Level: LOW
Final Vol: 5000 uL

Date Collected:
Date Received:
SDG No.: Q3757
Matrix: Water
% Solid: 0
Test: VOCMS Group3

CAS Number	Parameter	Conc.	Qua.	DF	MDL	LOQ / CRQL	Units	Date Ana.	BatchID
TARGETS									
75-01-4	Vinyl Chloride	0.26	U	1	0.26	1.00	ug/L	12/08/25 11:15	VX120825
75-35-4	1,1-Dichloroethene	0.23	U	1	0.23	1.00	ug/L	12/08/25 11:15	VX120825
75-34-3	1,1-Dichloroethane	0.23	U	1	0.23	1.00	ug/L	12/08/25 11:15	VX120825
156-59-2	cis-1,2-Dichloroethene	0.19	U	1	0.19	1.00	ug/L	12/08/25 11:15	VX120825
71-55-6	1,1,1-Trichloroethane	0.20	U	1	0.20	1.00	ug/L	12/08/25 11:15	VX120825
71-43-2	Benzene	0.15	U	1	0.15	1.00	ug/L	12/08/25 11:15	VX120825
107-06-2	1,2-Dichloroethane	0.22	U	1	0.22	1.00	ug/L	12/08/25 11:15	VX120825
79-01-6	Trichloroethene	0.090	U	1	0.090	1.00	ug/L	12/08/25 11:15	VX120825
79-00-5	1,1,2-Trichloroethane	0.21	U	1	0.21	1.00	ug/L	12/08/25 11:15	VX120825
127-18-4	Tetrachloroethene	0.23	U	1	0.23	1.00	ug/L	12/08/25 11:15	VX120825
SURROGATES									
17060-07-0	1,2-Dichloroethane-d4	49.5			70 (74) - 130 (125)	99%	SPK: 50		
1868-53-7	Dibromofluoromethane	46.0			70 (75) - 130 (124)	92%	SPK: 50		
2037-26-5	Toluene-d8	45.2			70 (86) - 130 (113)	90%	SPK: 50		
460-00-4	4-Bromofluorobenzene	55.8			70 (77) - 130 (121)	112%	SPK: 50		
INTERNAL STANDARDS									
		Area Count							
363-72-4	Pentafluorobenzene	173000							
540-36-3	1,4-Difluorobenzene	341000							
3114-55-4	Chlorobenzene-d5	370000							
3855-82-1	1,4-Dichlorobenzene-d4	180000							

U = Not Detected

LOQ = Limit of Quantitation

MDL = Method Detection Limit

LOD = Limit of Detection

E = Value Exceeds Calibration Range

Q = indicates LCS control criteria did not meet requirements

M = MS/MSD acceptance criteria did not meet requirements

J = Estimated Value

B = Analyte Found in Associated Method Blank

N = Presumptive Evidence of a Compound

* = Values outside of QC limits

D = Dilution

() = Laboratory InHouse Limit

A = Aldol-Condensation Reaction Products

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120825\
 Data File : VX048744.D
 Acq On : 08 Dec 2025 11:15
 Operator : JC/MD
 Sample : VX1208WBL01
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 VX1208WBL01

Quant Time: Dec 09 04:06:30 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X120425W.M
 Quant Title : SW846 8260
 QLast Update : Fri Dec 05 03:27:34 2025
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.531	168	173343	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.739	114	340953	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.037	117	370451	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.006	152	179749	50.000	ug/l	0.00

System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.934	65	115181	49.507	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	99.020%
35) Dibromofluoromethane	5.361	113	107991	46.001	ug/l	-0.01
Spiked Amount	50.000	Range	75 - 124	Recovery	=	92.000%
50) Toluene-d8	8.629	98	372131	45.222	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	90.440%#
62) 4-Bromofluorobenzene	11.061	95	158378	55.813	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	111.620%

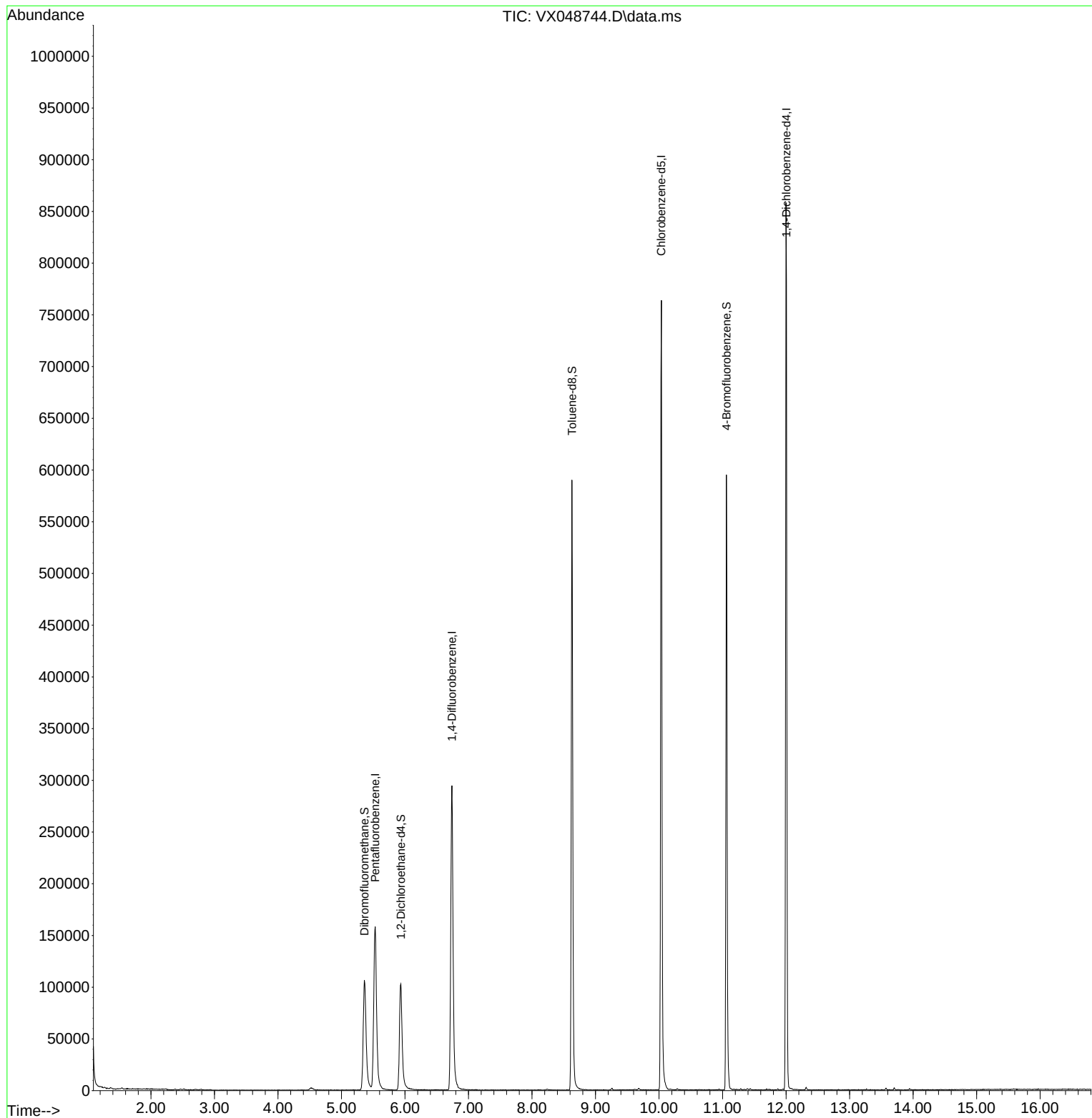
Target Compounds	Qvalue

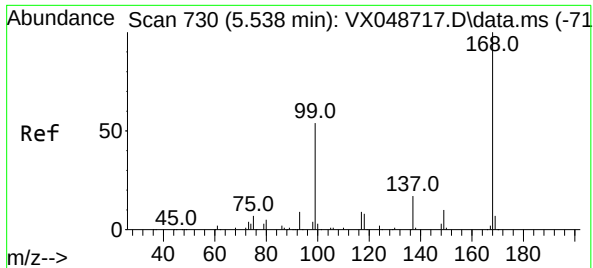
(#) = qualifier out of range (m) = manual integration (+) = signals summed

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120825\
Data File : VX048744.D
Acq On : 08 Dec 2025 11:15
Operator : JC/MD
Sample : VX1208WBL01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 4 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
VX1208WBL01

Quant Time: Dec 09 04:06:30 2025
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X120425W.M
Quant Title : SW846 8260
QLast Update : Fri Dec 05 03:27:34 2025
Response via : Initial Calibration

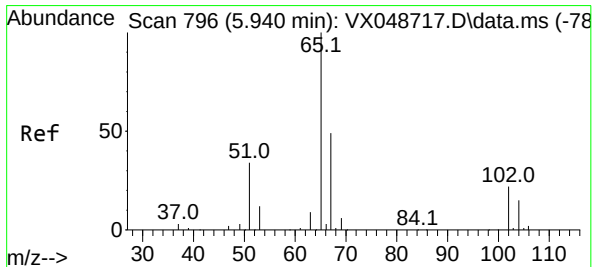
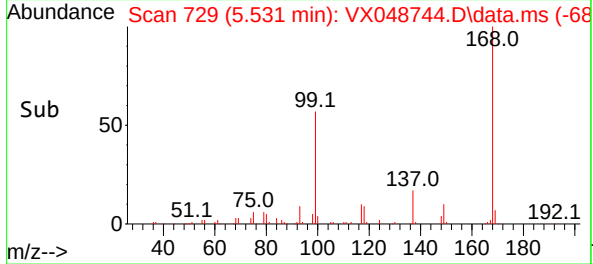
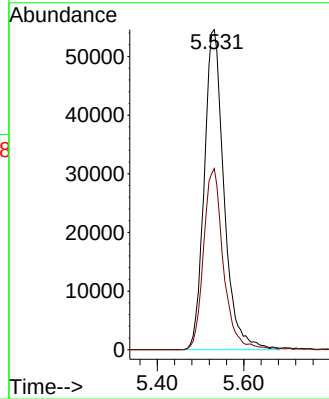
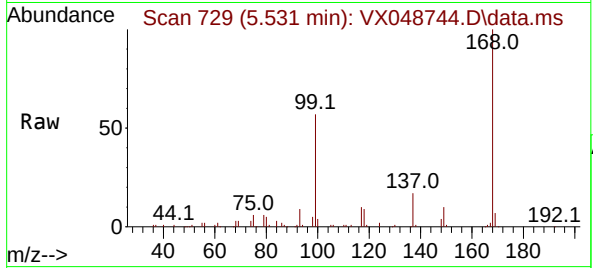




#1
Pentafluorobenzene
Concen: 50.000 ug/l
RT: 5.531 min Scan# 71
Delta R.T. -0.007 min
Lab File: VX048744.D
Acq: 08 Dec 2025 11:15

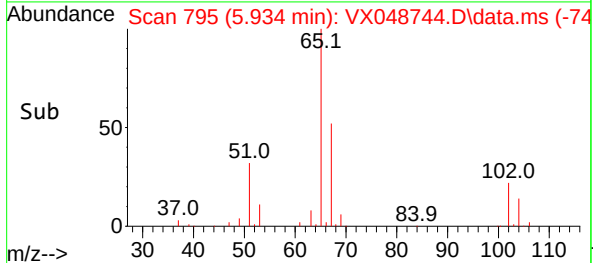
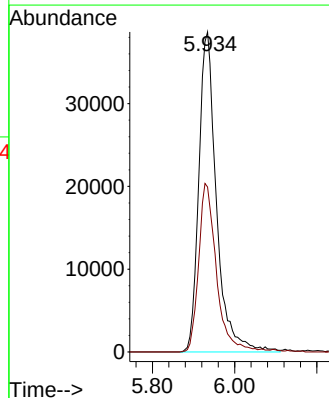
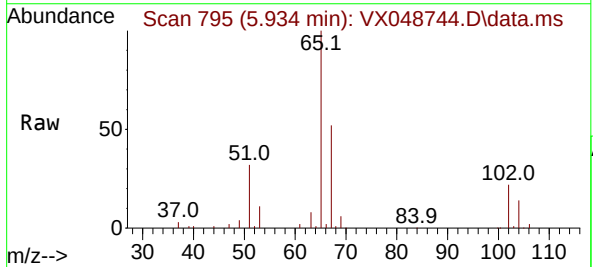
Instrument :
MSVOA_X
ClientSampleId :
VX1208WBL01

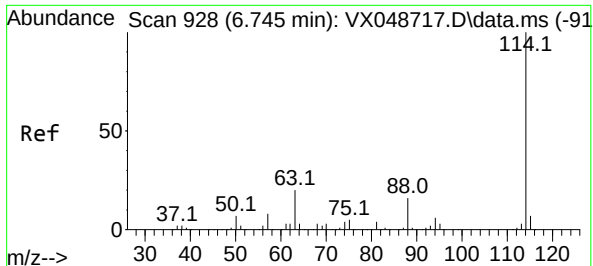
Tgt Ion:168 Resp: 173343
Ion Ratio Lower Upper
168 100
99 56.6 44.2 66.4



#33
1,2-Dichloroethane-d4
Concen: 49.507 ug/l
RT: 5.934 min Scan# 795
Delta R.T. -0.006 min
Lab File: VX048744.D
Acq: 08 Dec 2025 11:15

Tgt Ion: 65 Resp: 115181
Ion Ratio Lower Upper
65 100
67 52.8 0.0 107.4





#34

1,4-Difluorobenzene

Concen: 50.000 ug/l

RT: 6.739 min Scan# 91

Delta R.T. -0.006 min

Lab File: VX048744.D

Acq: 08 Dec 2025 11:15

Instrument :

MSVOA_X

ClientSampleId :

VX1208WBL01

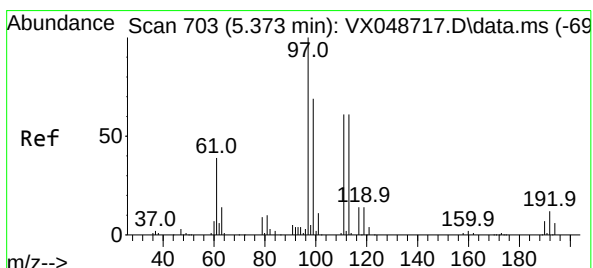
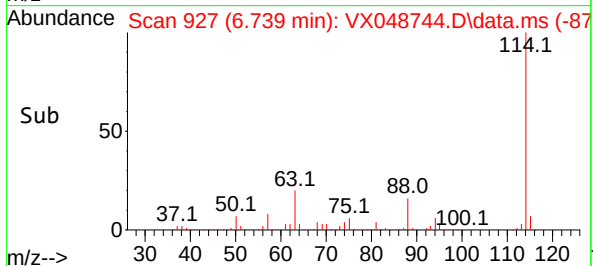
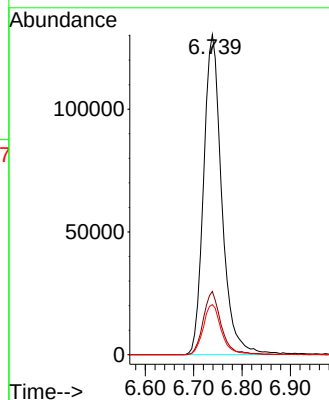
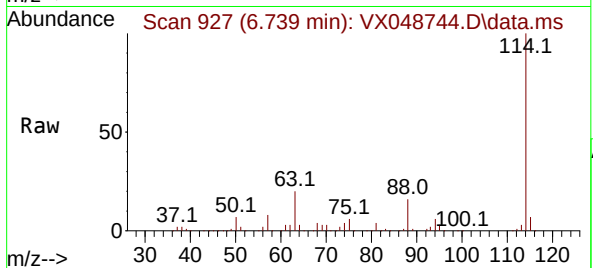
Tgt Ion:114 Resp: 340953

Ion Ratio Lower Upper

114 100

63 19.8 0.0 39.0

88 15.6 0.0 31.8



#35

Dibromofluoromethane

Concen: 46.001 ug/l

RT: 5.361 min Scan# 701

Delta R.T. -0.012 min

Lab File: VX048744.D

Acq: 08 Dec 2025 11:15

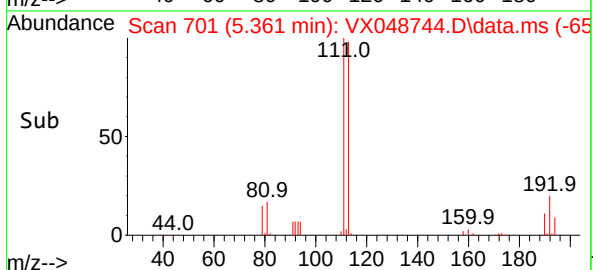
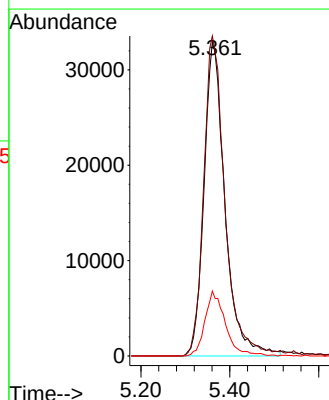
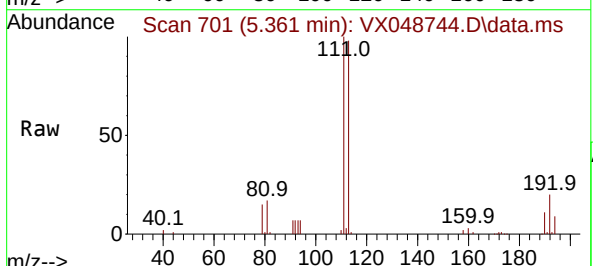
Tgt Ion:113 Resp: 107991

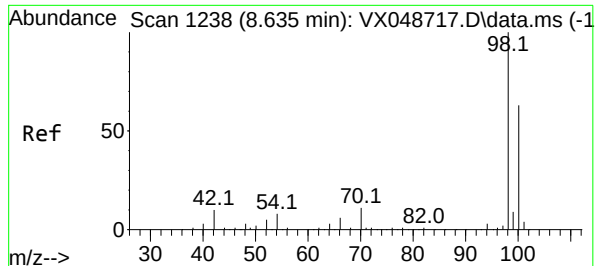
Ion Ratio Lower Upper

113 100

111 102.8 79.5 119.3

192 20.1 16.1 24.1





#50

Toluene-d8

Concen: 45.222 ug/l

RT: 8.629 min Scan# 11

Delta R.T. -0.006 min

Lab File: VX048744.D

Acq: 08 Dec 2025 11:15

Instrument :

MSVOA_X

ClientSampleId :

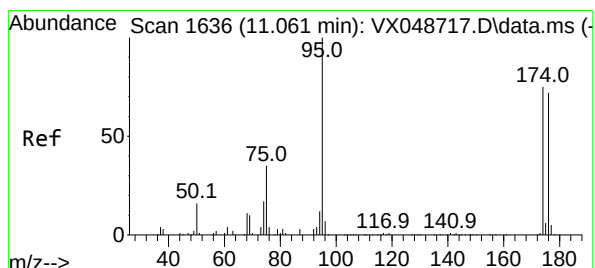
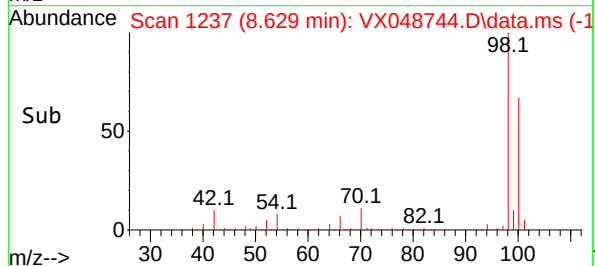
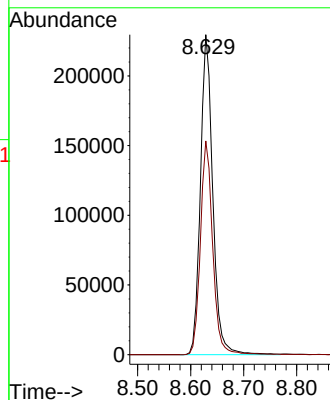
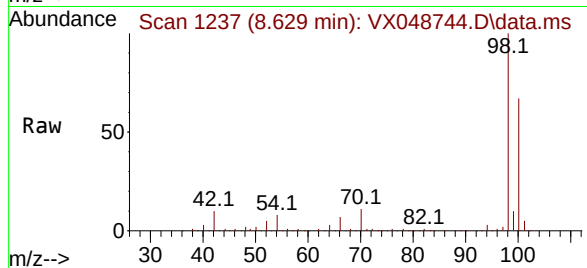
VX1208WBL01

Tgt Ion: 98 Resp: 372131

Ion Ratio Lower Upper

98 100

100 66.7 53.4 80.0



#62

4-Bromofluorobenzene

Concen: 55.813 ug/l

RT: 11.061 min Scan# 1636

Delta R.T. -0.000 min

Lab File: VX048744.D

Acq: 08 Dec 2025 11:15

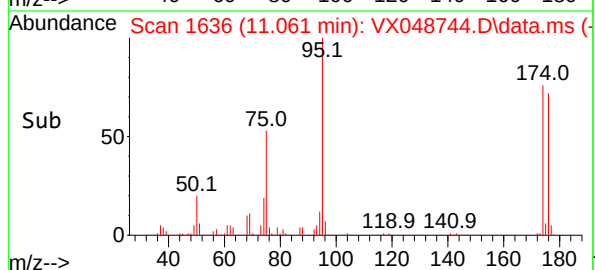
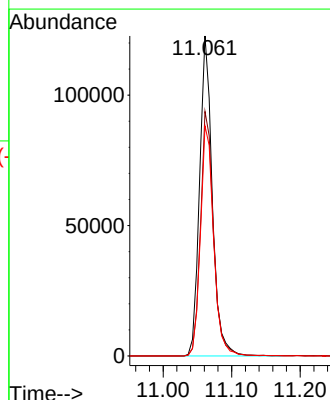
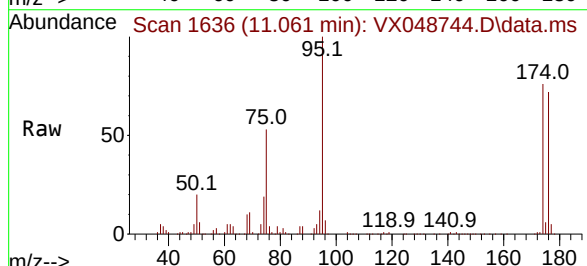
Tgt Ion: 95 Resp: 158378

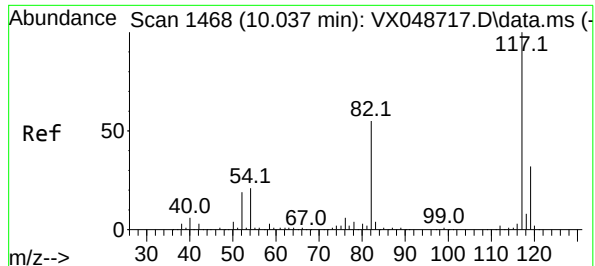
Ion Ratio Lower Upper

95 100

174 78.8 0.0 157.8

176 74.8 0.0 154.0





#63

Chlorobenzene-d5

Concen: 50.000 ug/l

RT: 10.037 min Scan# 1468

Delta R.T. -0.000 min

Lab File: VX048744.D

Acq: 08 Dec 2025 11:15

Instrument :

MSVOA_X

ClientSampleId :

VX1208WBL01

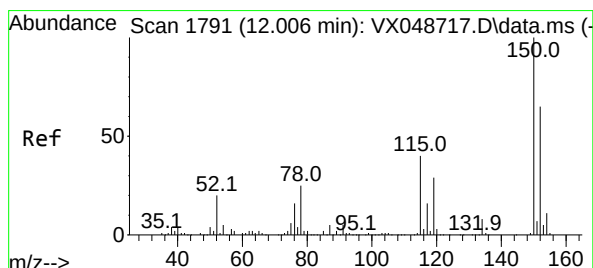
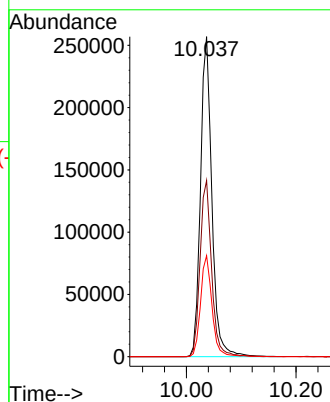
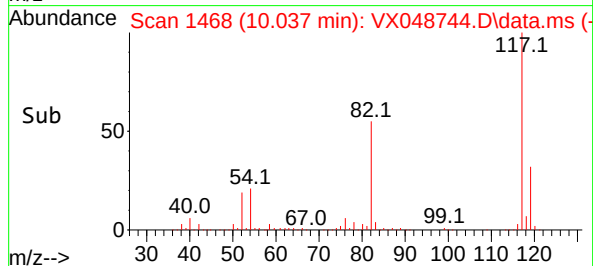
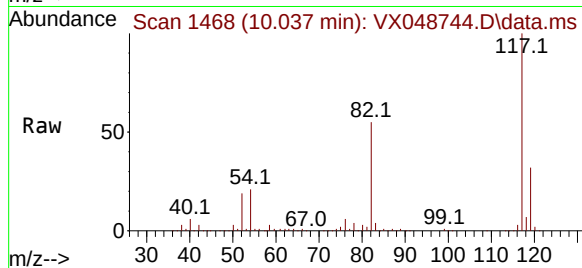
Tgt Ion:117 Resp: 370451

Ion Ratio Lower Upper

117 100

82 55.1 44.1 66.1

119 31.5 25.2 37.8



#72

1,4-Dichlorobenzene-d4

Concen: 50.000 ug/l

RT: 12.006 min Scan# 1791

Delta R.T. -0.000 min

Lab File: VX048744.D

Acq: 08 Dec 2025 11:15

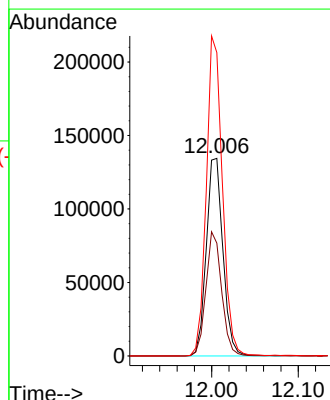
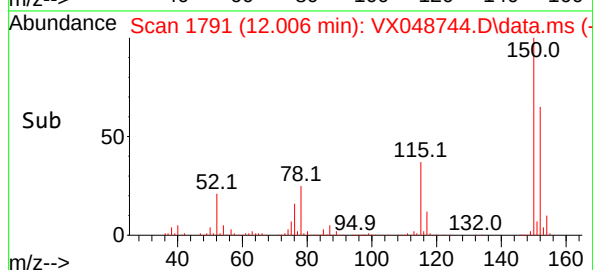
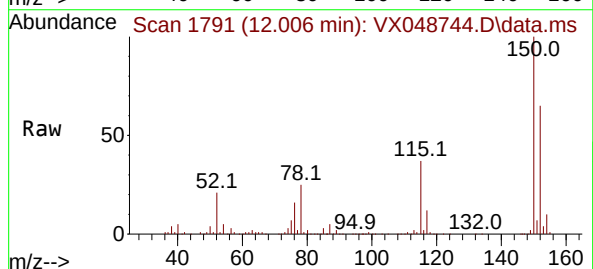
Tgt Ion:152 Resp: 179749

Ion Ratio Lower Upper

152 100

115 59.4 42.1 126.4

150 155.9 0.0 347.8





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

Report of Analysis

Client: JACOBS Engineering Group, Inc.
Project: Former Schlumberger STC PTC Site D3868221
Client Sample ID: VN1204WBS01
Lab Sample ID: VN1204WBS01
Analytical Method: 8260D
Sample Wt/Vol: 5 mL

Level: LOW
Final Vol: 5000 uL

Date Collected:
Date Received:
SDG No.: Q3757
Matrix: Water
% Solid: 0
Test: VOCMS Group3

CAS Number	Parameter	Conc.	Qua.	DF	MDL	LOQ / CRQL	Units	Date Ana.	BatchID
TARGETS									
75-01-4	Vinyl Chloride	20.3		1	0.26	1.00	ug/L	12/04/25 14:12	VN120425
75-35-4	1,1-Dichloroethene	20.4		1	0.23	1.00	ug/L	12/04/25 14:12	VN120425
75-34-3	1,1-Dichloroethane	19.6		1	0.23	1.00	ug/L	12/04/25 14:12	VN120425
156-59-2	cis-1,2-Dichloroethene	19.0		1	0.19	1.00	ug/L	12/04/25 14:12	VN120425
71-55-6	1,1,1-Trichloroethane	20.7		1	0.20	1.00	ug/L	12/04/25 14:12	VN120425
71-43-2	Benzene	19.3		1	0.15	1.00	ug/L	12/04/25 14:12	VN120425
107-06-2	1,2-Dichloroethane	20.8		1	0.22	1.00	ug/L	12/04/25 14:12	VN120425
79-01-6	Trichloroethene	18.3		1	0.090	1.00	ug/L	12/04/25 14:12	VN120425
79-00-5	1,1,2-Trichloroethane	19.9		1	0.21	1.00	ug/L	12/04/25 14:12	VN120425
127-18-4	Tetrachloroethene	18.1		1	0.23	1.00	ug/L	12/04/25 14:12	VN120425
SURROGATES									
17060-07-0	1,2-Dichloroethane-d4	53.8			70 (74) - 130 (125)	108%	SPK: 50		
1868-53-7	Dibromofluoromethane	50.7			70 (75) - 130 (124)	101%	SPK: 50		
2037-26-5	Toluene-d8	51.7			70 (86) - 130 (113)	103%	SPK: 50		
460-00-4	4-Bromofluorobenzene	54.4			70 (77) - 130 (121)	109%	SPK: 50		
INTERNAL STANDARDS									
		Area Count							
363-72-4	Pentafluorobenzene	304000							
540-36-3	1,4-Difluorobenzene	560000							
3114-55-4	Chlorobenzene-d5	492000							
3855-82-1	1,4-Dichlorobenzene-d4	238000							

U = Not Detected

LOQ = Limit of Quantitation

MDL = Method Detection Limit

LOD = Limit of Detection

E = Value Exceeds Calibration Range

Q = indicates LCS control criteria did not meet requirements

M = MS/MSD acceptance criteria did not meet requirements

J = Estimated Value

B = Analyte Found in Associated Method Blank

N = Presumptive Evidence of a Compound

* = Values outside of QC limits

D = Dilution

() = Laboratory InHouse Limit

A = Aldol-Condensation Reaction Products

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN120425\
 Data File : VN088443.D
 Acq On : 04 Dec 2025 14:12
 Operator : JC\MD
 Sample : VN1204WBS01
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 VN1204WBS01

Manual Integrations
 APPROVED

Reviewed By :John Carlone 12/05/2025
 Supervised By :Semsettin Yesilyurt 12/05/2025

Quant Time: Dec 05 00:14:59 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N112425W.M
 Quant Title : SW846 8260
 QLast Update : Tue Nov 25 01:34:52 2025
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.206	168	303745	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.077	114	559564	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.841	117	491842	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.770	152	237731	50.000	ug/l	0.00

System Monitoring Compounds

33) 1,2-Dichloroethane-d4	8.559	65	307567	53.752	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	= 107.500%	
35) Dibromofluoromethane	8.147	113	196606	50.701	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 101.400%	
50) Toluene-d8	10.547	98	745425	51.664	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	= 103.320%	
62) 4-Bromofluorobenzene	12.823	95	269159	54.369	ug/l	0.00
Spiked Amount	50.000	Range	77 - 121	Recovery	= 108.740%	

Target Compounds

						Qvalue
2) Dichlorodifluoromethane	2.142	85	97797	21.436	ug/l	93
3) Chloromethane	2.383	50	93145	19.009	ug/l	96
4) Vinyl Chloride	2.536	62	100990	20.259	ug/l	100
5) Bromomethane	2.983	94	46454	19.362	ug/l	91
6) Chloroethane	3.142	64	62589	20.289	ug/l	98
7) Trichlorofluoromethane	3.512	101	150074	21.421	ug/l	95
8) Diethyl Ether	3.959	74	53546	20.363	ug/l	99
9) 1,1,2-Trichlorotrifluo...	4.371	101	78946	21.166	ug/l	99
10) Methyl Iodide	4.583	142	63657	17.241	ug/l	93
11) Tert butyl alcohol	5.506	59	76083	83.283	ug/l	97
12) 1,1-Dichloroethene	4.342	96	75627	20.408	ug/l	96
13) Acrolein	4.183	56	95364	109.036	ug/l	98
14) Allyl chloride	5.012	41	131363	19.125	ug/l	97
15) Acrylonitrile	5.706	53	244233	92.662	ug/l	99
16) Acetone	4.424	43	183497	88.674	ug/l	96
17) Carbon Disulfide	4.712	76	223379	18.447	ug/l	96
18) Methyl Acetate	5.018	43	120530	19.567	ug/l	99
19) Methyl tert-butyl Ether	5.783	73	272930	19.658	ug/l	99
20) Methylene Chloride	5.265	84	81214	18.343	ug/l	99
21) trans-1,2-Dichloroethene	5.777	96	79259	19.308	ug/l	98
22) Diisopropyl ether	6.659	45	293315	20.002	ug/l	98
23) Vinyl Acetate	6.588	43	1278749	108.235	ug/l	99
24) 1,1-Dichloroethane	6.553	63	162999	19.628	ug/l	97
25) 2-Butanone	7.465	43	310201	93.245	ug/l	100
26) 2,2-Dichloropropane	7.471	77	150702	21.025	ug/l	97
27) cis-1,2-Dichloroethene	7.471	96	90646	19.000	ug/l	96
28) Bromochloromethane	7.794	49	70911	17.142	ug/l	98
29) Tetrahydrofuran	7.818	42	228135	95.607	ug/l	98
30) Chloroform	7.953	83	165040	20.226	ug/l	100
31) Cyclohexane	8.241	56	151560	20.140	ug/l	99
32) 1,1,1-Trichloroethane	8.147	97	148393	20.683	ug/l	97
36) 1,1-Dichloropropene	8.353	75	116928	19.538	ug/l	98
37) Ethyl Acetate	7.547	43	137369	17.457	ug/l	97
38) Carbon Tetrachloride	8.341	117	123356	20.706	ug/l	98
39) Methylcyclohexane	9.582	83	125506	19.606	ug/l	90
40) Benzene	8.582	78	339466	19.326	ug/l	99

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN120425\
 Data File : VN088443.D
 Acq On : 04 Dec 2025 14:12
 Operator : JC\MD
 Sample : VN1204WBS01
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 VN1204WBS01

Manual Integrations
 APPROVED

Reviewed By :John Carlone 12/05/2025
 Supervised By :Semsettin Yesilyurt 12/05/2025

Quant Time: Dec 05 00:14:59 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N112425W.M
 Quant Title : SW846 8260
 QLast Update : Tue Nov 25 01:34:52 2025
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Methacrylonitrile	7.765	41	74031	18.995	ug/l	99
42) 1,2-Dichloroethane	8.647	62	139506	20.817	ug/l	97
43) Isopropyl Acetate	8.671	43	224127	18.642	ug/l	99
44) Trichloroethene	9.329	130	76077	18.330	ug/l	97
45) 1,2-Dichloropropane	9.600	63	87062	19.346	ug/l	99
46) Dibromomethane	9.688	93	62387	19.425	ug/l	99
47) Bromodichloromethane	9.865	83	130611	19.931	ug/l	99
48) Methyl methacrylate	9.659	41	104044	18.966	ug/l	99
49) 1,4-Dioxane	9.677	88	21834	312.972	ug/l #	94
51) 4-Methyl-2-Pentanone	10.424	43	704216	99.372	ug/l	100
52) Toluene	10.606	92	205014	20.214	ug/l	99
53) t-1,3-Dichloropropene	10.812	75	131896	19.609	ug/l	100
54) cis-1,3-Dichloropropene	10.288	75	137564	19.447	ug/l	96
55) 1,1,2-Trichloroethane	10.994	97	77993	19.876	ug/l	96
56) Ethyl methacrylate	10.853	69	122489	19.761	ug/l	99
57) 1,3-Dichloropropane	11.141	76	141160	19.940	ug/l	99
58) 2-Chloroethyl Vinyl ether	10.135	63	316457	88.556	ug/l	98
59) 2-Hexanone	11.176	43	422202	94.573	ug/l	99
60) Dibromochloromethane	11.335	129	89380	19.980	ug/l	100
61) 1,2-Dibromoethane	11.447	107	78746	19.817	ug/l	98
64) Tetrachloroethene	11.082	164	70474	18.074	ug/l	94
65) Chlorobenzene	11.870	112	218454	19.774	ug/l	100
66) 1,1,1,2-Tetrachloroethane	11.941	131	72384	19.426	ug/l	96
67) Ethyl Benzene	11.941	91	390956	20.247	ug/l	98
68) m/p-Xylenes	12.047	106	286124	39.966	ug/l	96
69) o-Xylene	12.370	106	139831	20.487	ug/l	99
70) Styrene	12.388	104	232254	20.957	ug/l	99
71) Bromoform	12.553	173	55870	20.471	ug/l #	99
73) Isopropylbenzene	12.670	105	360370	20.502	ug/l	99
74) N-amyl acetate	12.688	43	116340m	19.435	ug/l	
75) 1,1,2,2-Tetrachloroethane	12.912	83	112119	19.731	ug/l	99
76) 1,2,3-Trichloropropane	12.970	75	113577m	20.398	ug/l	
77) Bromobenzene	12.953	156	82378	20.483	ug/l	99
78) n-propylbenzene	13.012	91	452912	21.062	ug/l	99
79) 2-Chlorotoluene	13.100	91	273041	20.792	ug/l	98
80) 1,3,5-Trimethylbenzene	13.147	105	301838	20.723	ug/l	100
81) trans-1,4-Dichloro-2-b...	12.723	75	35744	18.212	ug/l #	91
82) 4-Chlorotoluene	13.194	91	286032	21.930	ug/l	97
83) tert-Butylbenzene	13.412	119	261764	21.141	ug/l	97
84) 1,2,4-Trimethylbenzene	13.459	105	312793	21.606	ug/l	97
85) sec-Butylbenzene	13.594	105	375613	21.117	ug/l	99
86) p-Isopropyltoluene	13.706	119	306028	20.981	ug/l	97
87) 1,3-Dichlorobenzene	13.712	146	154606	19.748	ug/l	99
88) 1,4-Dichlorobenzene	13.788	146	162379	19.800	ug/l	99
89) n-Butylbenzene	14.029	91	297476	21.005	ug/l	98
90) Hexachloroethane	14.306	117	52252	19.610	ug/l	98
91) 1,2-Dichlorobenzene	14.082	146	147868	20.071	ug/l	98
92) 1,2-Dibromo-3-Chloropr...	14.694	75	28137	20.408	ug/l	99
93) 1,2,4-Trichlorobenzene	15.364	180	83932	19.174	ug/l	95
94) Hexachlorobutadiene	15.470	225	28512	18.126	ug/l	97
95) Naphthalene	15.611	128	307489	20.271	ug/l	100
96) 1,2,3-Trichlorobenzene	15.811	180	83257	19.543	ug/l	99

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN120425\
Data File : VN088443.D
Acq On : 04 Dec 2025 14:12
Operator : JC\MD
Sample : VN1204WBS01
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
VN1204WBS01

Manual Integrations
APPROVED

Reviewed By :John Carlone 12/05/2025
Supervised By :Semsettin Yesilyurt 12/05/2025

Quant Time: Dec 05 00:14:59 2025
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N112425W.M
Quant Title : SW846 8260
QLast Update : Tue Nov 25 01:34:52 2025
Response via : Initial Calibration

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

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

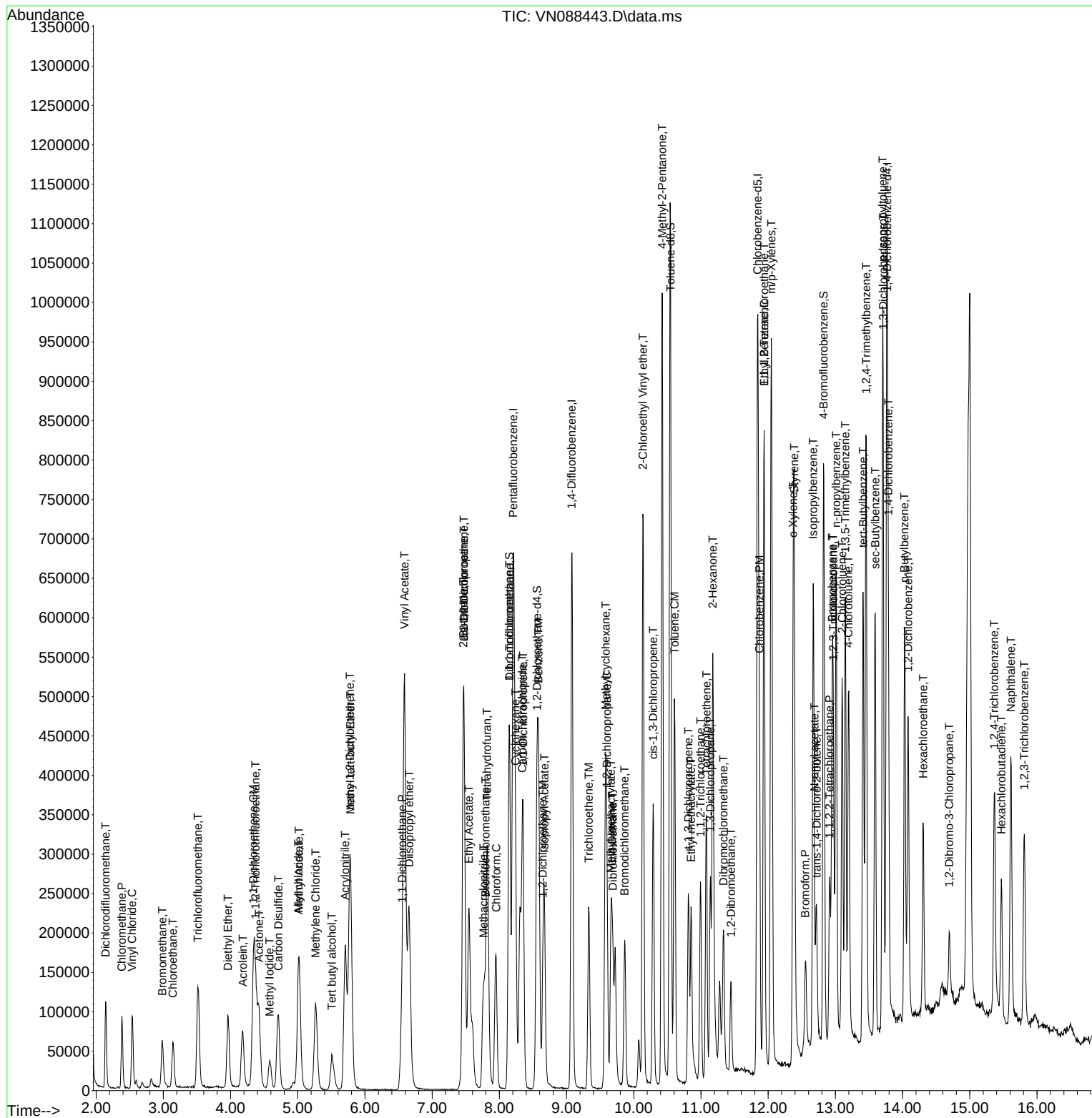
```
Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN120425\  
Data File : VN088443.D  
Acq On    : 04 Dec 2025   14:12  
Operator  : JC\MD  
Sample    : VN1204WBS01  
Misc      : 5.0mL/MSVOA_N/WATER  
ALS Vial  : 9      Sample Multiplier: 1
```

Instrument :
MSVOA_N
ClientSampleId :
VN1204WBS01

Manual Integrations

Reviewed By :John Carlone 12/05/2025
Supervised By :Semsettin Yesilyurt 12/05/2025

Quant Time: Dec 05 00:14:59 2025
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N112425W.M
Quant Title : SW846 8260
QLast Update : Tue Nov 25 01:34:52 2025
Response via : Initial Calibration





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

Report of Analysis

Client: JACOBS Engineering Group, Inc.
Project: Former Schlumberger STC PTC Site D3868221
Client Sample ID: VX1205WBS01
Lab Sample ID: VX1205WBS01
Analytical Method: 8260D
Sample Wt/Vol: 5 mL

Level: LOW
Final Vol: 5000 uL

Date Collected:
Date Received:
SDG No.: Q3757
Matrix: Water
% Solid: 0
Test: VOCMS Group3

CAS Number	Parameter	Conc.	Qua.	DF	MDL	LOQ / CRQL	Units	Date Ana.	BatchID
TARGETS									
75-01-4	Vinyl Chloride	16.1		1	0.26	1.00	ug/L	12/05/25 11:31	VX120525
75-35-4	1,1-Dichloroethene	19.4		1	0.23	1.00	ug/L	12/05/25 11:31	VX120525
75-34-3	1,1-Dichloroethane	20.1		1	0.23	1.00	ug/L	12/05/25 11:31	VX120525
156-59-2	cis-1,2-Dichloroethene	19.1		1	0.19	1.00	ug/L	12/05/25 11:31	VX120525
71-55-6	1,1,1-Trichloroethane	20.0		1	0.20	1.00	ug/L	12/05/25 11:31	VX120525
71-43-2	Benzene	18.7		1	0.15	1.00	ug/L	12/05/25 11:31	VX120525
107-06-2	1,2-Dichloroethane	20.1		1	0.22	1.00	ug/L	12/05/25 11:31	VX120525
79-01-6	Trichloroethene	19.0		1	0.090	1.00	ug/L	12/05/25 11:31	VX120525
79-00-5	1,1,2-Trichloroethane	19.1		1	0.21	1.00	ug/L	12/05/25 11:31	VX120525
127-18-4	Tetrachloroethene	17.6		1	0.23	1.00	ug/L	12/05/25 11:31	VX120525
SURROGATES									
17060-07-0	1,2-Dichloroethane-d4	52.4			70 (74) - 130 (125)	105%	SPK: 50		
1868-53-7	Dibromofluoromethane	48.0			70 (75) - 130 (124)	96%	SPK: 50		
2037-26-5	Toluene-d8	49.6			70 (86) - 130 (113)	99%	SPK: 50		
460-00-4	4-Bromofluorobenzene	52.5			70 (77) - 130 (121)	105%	SPK: 50		
INTERNAL STANDARDS									
		Area Count							
363-72-4	Pentafluorobenzene	179000							
540-36-3	1,4-Difluorobenzene	294000							
3114-55-4	Chlorobenzene-d5	269000							
3855-82-1	1,4-Dichlorobenzene-d4	133000							

U = Not Detected

LOQ = Limit of Quantitation

MDL = Method Detection Limit

LOD = Limit of Detection

E = Value Exceeds Calibration Range

Q = indicates LCS control criteria did not meet requirements

M = MS/MSD acceptance criteria did not meet requirements

J = Estimated Value

B = Analyte Found in Associated Method Blank

N = Presumptive Evidence of a Compound

* = Values outside of QC limits

D = Dilution

() = Laboratory InHouse Limit

A = Aldol-Condensation Reaction Products

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120525\
 Data File : VX048726.D
 Acq On : 05 Dec 2025 11:31
 Operator : JC/MD
 Sample : VX1205WBS01
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 VX1205WBS01

Manual Integrations
 APPROVED

Reviewed By : John Carlone 12/08/2025
 Supervised By : Semsettin Yesilyurt 12/08/2025

Quant Time: Dec 05 21:59:38 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X120425W.M
 Quant Title : SW846 8260
 QLast Update : Fri Dec 05 03:27:34 2025
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.525	168	179194	50.000	ug/l	-0.01
34) 1,4-Difluorobenzene	6.739	114	293931	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.037	117	269397	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.006	152	132992	50.000	ug/l	0.00

System Monitoring Compounds

33) 1,2-Dichloroethane-d4	5.934	65	125979	52.380	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	= 104.760%	
35) Dibromofluoromethane	5.361	113	97103	47.980	ug/l	-0.01
Spiked Amount	50.000	Range	75 - 124	Recovery	= 95.960%	
50) Toluene-d8	8.629	98	352069	49.628	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	= 99.260%	
62) 4-Bromofluorobenzene	11.061	95	128406	52.490	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 104.980%	

Target Compounds

						Qvalue
2) Dichlorodifluoromethane	1.173	85	41863	22.929	ug/l	96
3) Chloromethane	1.301	50	44995	19.530	ug/l	97
4) Vinyl Chloride	1.380	62	39598	16.138	ug/l	97
5) Bromomethane	1.618	94	28574	18.844	ug/l	97
6) Chloroethane	1.697	64	24297	15.910	ug/l	92
7) Trichlorofluoromethane	1.898	101	68676	19.528	ug/l	97
8) Diethyl Ether	2.130	74	23437	16.697	ug/l	97
9) 1,1,2-Trichlorotrifluo...	2.337	101	43147	21.164	ug/l	99
10) Methyl Iodide	2.453	142	61205	21.087	ug/l	99
11) Tert butyl alcohol	2.928	59	35673	91.869	ug/l	100
12) 1,1-Dichloroethene	2.325	96	40443	19.436	ug/l	99
13) Acrolein	2.233	56	31115	109.833	ug/l	100
14) Allyl chloride	2.660	41	71931	20.590	ug/l	94
15) Acrylonitrile	3.050	53	110416	91.258	ug/l	99
16) Acetone	2.367	43	87721	95.274	ug/l	97
17) Carbon Disulfide	2.514	76	125337	20.742	ug/l	100
18) Methyl Acetate	2.697	43	47862	18.476	ug/l	99
19) Methyl tert-butyl Ether	3.105	73	128879	19.031	ug/l	97
20) Methylene Chloride	2.782	84	41148	17.880	ug/l	95
21) trans-1,2-Dichloroethene	3.087	96	40211	18.934	ug/l	93
22) Diisopropyl ether	3.739	45	131789	20.185	ug/l #	80
23) Vinyl Acetate	3.709	43	567718	103.483	ug/l	97
24) 1,1-Dichloroethane	3.599	63	73078	20.071	ug/l	100
25) 2-Butanone	4.526	43	144944	106.663	ug/l	91
26) 2,2-Dichloropropane	4.459	77	65233	20.249	ug/l	98
27) cis-1,2-Dichloroethene	4.471	96	45731	19.082	ug/l	92
28) Bromochloromethane	4.879	49	32271	18.083	ug/l	94
29) Tetrahydrofuran	4.977	42	104631	100.837	ug/l	95
30) Chloroform	5.068	83	77868	19.960	ug/l	99
31) Cyclohexane	5.452	56	62885	19.345	ug/l	95
32) 1,1,1-Trichloroethane	5.361	97	69932	20.006	ug/l	96
36) 1,1-Dichloropropene	5.672	75	57482	20.737	ug/l	100
37) Ethyl Acetate	4.684	43	65510	19.779	ug/l	99
38) Carbon Tetrachloride	5.653	117	65362	20.817	ug/l	94
39) Methylcyclohexane	7.360	83	62448	19.052	ug/l	97
40) Benzene	6.019	78	153807	18.718	ug/l	99

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120525\
 Data File : VX048726.D
 Acq On : 05 Dec 2025 11:31
 Operator : JC/MD
 Sample : VX1205WBS01
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 VX1205WBS01

Manual Integrations
 APPROVED

Reviewed By :John Carlone 12/08/2025
 Supervised By :Semsettin Yesilyurt 12/08/2025

Quant Time: Dec 05 21:59:38 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X120425W.M
 Quant Title : SW846 8260
 QLast Update : Fri Dec 05 03:27:34 2025
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Methacrylonitrile	4.897	41	29896	19.838	ug/l	97
42) 1,2-Dichloroethane	6.068	62	58414	20.085	ug/l	95
43) Isopropyl Acetate	6.312	43	97581	19.456	ug/l	95
44) Trichloroethene	7.104	130	40756	19.009	ug/l	98
45) 1,2-Dichloropropane	7.409	63	37372	18.196	ug/l	99
46) Dibromomethane	7.562	93	29281	18.800	ug/l	99
47) Bromodichloromethane	7.805	83	61760	19.019	ug/l	99
48) Methyl methacrylate	7.671	41	46528	18.442	ug/l	94
49) 1,4-Dioxane	7.641	88	11029	317.857	ug/l	96
51) 4-Methyl-2-Pentanone	8.549	43	315106	101.221	ug/l	96
52) Toluene	8.702	92	98680	19.915	ug/l	100
53) t-1,3-Dichloropropene	8.958	75	61457	19.894	ug/l	97
54) cis-1,3-Dichloropropene	8.348	75	63144	19.019	ug/l	91
55) 1,1,2-Trichloroethane	9.134	97	37998	19.059	ug/l	98
56) Ethyl methacrylate	9.098	69	60402	19.356	ug/l #	94
57) 1,3-Dichloropropane	9.293	76	68113	20.546	ug/l	100
58) 2-Chloroethyl Vinyl ether	8.220	63	144206	85.498	ug/l	99
59) 2-Hexanone	9.409	43	255149	116.758	ug/l	98
60) Dibromochloromethane	9.500	129	54125	21.955	ug/l	98
61) 1,2-Dibromoethane	9.592	107	47896	22.197	ug/l	100
64) Tetrachloroethene	9.256	164	34511	17.575	ug/l	97
65) Chlorobenzene	10.061	112	109860	18.630	ug/l	95
66) 1,1,1,2-Tetrachloroethane	10.147	131	37880	17.960	ug/l	99
67) Ethyl Benzene	10.177	91	190737	19.822	ug/l	97
68) m/p-Xylenes	10.281	106	154292	41.164	ug/l	100
69) o-Xylene	10.622	106	74050	21.083	ug/l	100
70) Styrene	10.634	104	126101	20.914	ug/l	99
71) Bromoform	10.781	173	34490	18.317	ug/l #	99
73) Isopropylbenzene	10.945	105	189027	21.092	ug/l	99
74) N-amyl acetate	10.823	43	88194	22.648	ug/l	99
75) 1,1,2,2-Tetrachloroethane	11.195	83	58063	19.201	ug/l	100
76) 1,2,3-Trichloropropane	11.220	75	56826m	22.941	ug/l	
77) Bromobenzene	11.177	156	43921	18.361	ug/l	99
78) n-propylbenzene	11.287	91	218146	21.016	ug/l	100
79) 2-Chlorotoluene	11.348	91	129011	20.383	ug/l	99
80) 1,3,5-Trimethylbenzene	11.433	105	150128	20.379	ug/l	98
81) trans-1,4-Dichloro-2-b...	11.000	75	20029	19.207	ug/l	94
82) 4-Chlorotoluene	11.433	91	153143	20.734	ug/l	99
83) tert-Butylbenzene	11.695	119	149901	19.584	ug/l	97
84) 1,2,4-Trimethylbenzene	11.732	105	152255	20.489	ug/l	100
85) sec-Butylbenzene	11.872	105	189342	20.411	ug/l	99
86) p-Isopropyltoluene	11.988	119	158652	20.008	ug/l	99
87) 1,3-Dichlorobenzene	11.951	146	84489	19.160	ug/l	98
88) 1,4-Dichlorobenzene	12.024	146	84209	18.479	ug/l	97
89) n-Butylbenzene	12.311	91	147367	19.849	ug/l	100
90) Hexachloroethane	12.518	117	30361	18.760	ug/l	98
91) 1,2-Dichlorobenzene	12.317	146	78982	18.475	ug/l	100
92) 1,2-Dibromo-3-Chloropr...	12.920	75	13457	18.285	ug/l	91
93) 1,2,4-Trichlorobenzene	13.567	180	48161	17.057	ug/l	99
94) Hexachlorobutadiene	13.707	225	21240	18.186	ug/l	95
95) Naphthalene	13.756	128	154818	20.946	ug/l	100
96) 1,2,3-Trichlorobenzene	13.939	180	45023	17.290	ug/l	97

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120525\
Data File : VX048726.D
Acq On : 05 Dec 2025 11:31
Operator : JC/MD
Sample : VX1205WBS01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 5 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
VX1205WBS01

Manual Integrations
APPROVED

Reviewed By :John Carlone 12/08/2025
Supervised By :Semsettin Yesilyurt 12/08/2025

Quant Time: Dec 05 21:59:38 2025
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X120425W.M
Quant Title : SW846 8260
QLast Update : Fri Dec 05 03:27:34 2025
Response via : Initial Calibration

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

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

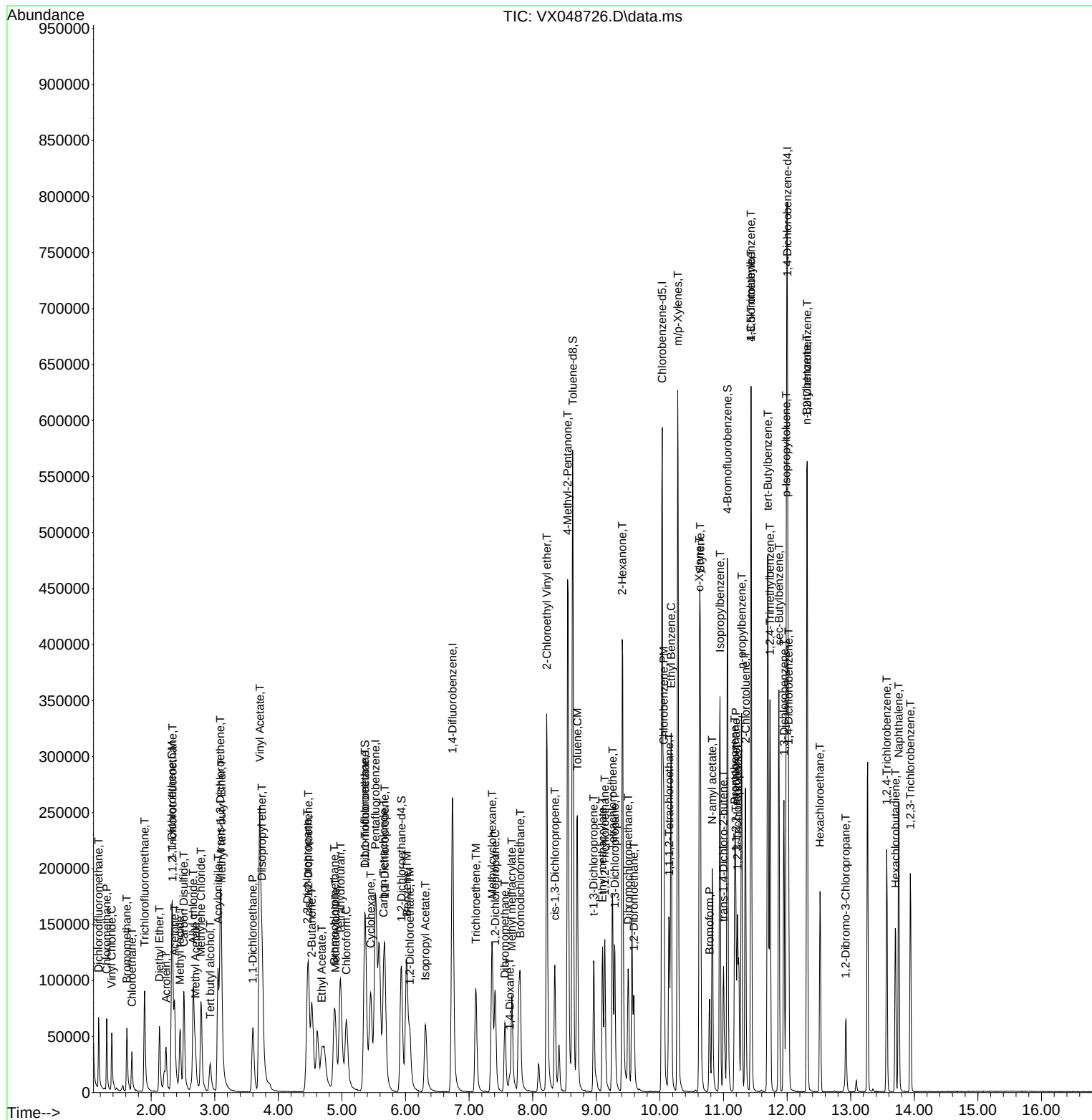

```
Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120525\
Data File : VX048726.D
Acq On    : 05 Dec 2025  11:31
Operator  : JC/MD
Sample    : VX1205WBS01
Misc      : 5.0mL/MSVOA_X/WATER
ALS Vial  : 5      Sample Multiplier: 1
```

Instrument :
MSVOA_X
ClientSampleId :
VX1205WBS01

Manual Integrations APPROVED

Reviewed By :John Carlone 12/08/2025
Supervised By :Semsettin Yesilyurt 12/08/2025

Quant Time: Dec 05 21:59:38 2025
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X120425W.M
Quant Title : SW846 8260
QLast Update : Fri Dec 05 03:27:34 2025
Response via : Initial Calibration





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

Report of Analysis

Client: JACOBS Engineering Group, Inc.
Project: Former Schlumberger STC PTC Site D3868221
Client Sample ID: VX1208WBS01
Lab Sample ID: VX1208WBS01
Analytical Method: 8260D
Sample Wt/Vol: 5 mL

Level: LOW
Final Vol: 5000 uL

Date Collected:
Date Received:
SDG No.: Q3757
Matrix: Water
% Solid: 0
Test: VOCMS Group3

CAS Number	Parameter	Conc.	Qua.	DF	MDL	LOQ / CRQL	Units	Date Ana.	BatchID
TARGETS									
75-01-4	Vinyl Chloride	19.8		1	0.26	1.00	ug/L	12/08/25 11:44	VX120825
75-35-4	1,1-Dichloroethene	19.2		1	0.23	1.00	ug/L	12/08/25 11:44	VX120825
75-34-3	1,1-Dichloroethane	20.8		1	0.23	1.00	ug/L	12/08/25 11:44	VX120825
156-59-2	cis-1,2-Dichloroethene	19.9		1	0.19	1.00	ug/L	12/08/25 11:44	VX120825
71-55-6	1,1,1-Trichloroethane	20.1		1	0.20	1.00	ug/L	12/08/25 11:44	VX120825
71-43-2	Benzene	18.8		1	0.15	1.00	ug/L	12/08/25 11:44	VX120825
107-06-2	1,2-Dichloroethane	21.4		1	0.22	1.00	ug/L	12/08/25 11:44	VX120825
79-01-6	Trichloroethene	17.6		1	0.090	1.00	ug/L	12/08/25 11:44	VX120825
79-00-5	1,1,2-Trichloroethane	21.7		1	0.21	1.00	ug/L	12/08/25 11:44	VX120825
127-18-4	Tetrachloroethene	16.5		1	0.23	1.00	ug/L	12/08/25 11:44	VX120825
SURROGATES									
17060-07-0	1,2-Dichloroethane-d4	56.7			70 (74) - 130 (125)	113%	SPK: 50		
1868-53-7	Dibromofluoromethane	50.4			70 (75) - 130 (124)	101%	SPK: 50		
2037-26-5	Toluene-d8	50.6			70 (86) - 130 (113)	101%	SPK: 50		
460-00-4	4-Bromofluorobenzene	58.8			70 (77) - 130 (121)	118%	SPK: 50		
INTERNAL STANDARDS									
		Area Count							
363-72-4	Pentafluorobenzene	178000							
540-36-3	1,4-Difluorobenzene	299000							
3114-55-4	Chlorobenzene-d5	311000							
3855-82-1	1,4-Dichlorobenzene-d4	158000							

U = Not Detected

LOQ = Limit of Quantitation

MDL = Method Detection Limit

LOD = Limit of Detection

E = Value Exceeds Calibration Range

Q = indicates LCS control criteria did not meet requirements

M = MS/MSD acceptance criteria did not meet requirements

J = Estimated Value

B = Analyte Found in Associated Method Blank

N = Presumptive Evidence of a Compound

* = Values outside of QC limits

D = Dilution

() = Laboratory InHouse Limit

A = Aldol-Condensation Reaction Products

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120825\
 Data File : VX048745.D
 Acq On : 08 Dec 2025 11:44
 Operator : JC/MD
 Sample : VX1208WBS01
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 VX1208WBS01

Manual Integrations
 APPROVED

Reviewed By :John Carlone 12/09/2025
 Supervised By :Mahesh Dadoda 12/09/2025

Quant Time: Dec 09 04:06:52 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X120425W.M
 Quant Title : SW846 8260
 QLast Update : Fri Dec 05 03:27:34 2025
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.531	168	178446	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.738	114	298973	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.037	117	311271	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.000	152	157846	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.928	65	135832	56.713	ug/l	-0.01
Spiked Amount 50.000	Range 78 - 117		Recovery = 113.420%			
35) Dibromofluoromethane	5.367	113	103655	50.354	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery = 100.700%			
50) Toluene-d8	8.628	98	364834	50.560	ug/l	0.00
Spiked Amount 50.000	Range 92 - 112		Recovery = 101.120%			
62) 4-Bromofluorobenzene	11.061	95	146293	58.794	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery = 117.580%			
Target Compounds					Qvalue	
2) Dichlorodifluoromethane	1.172	85	41301	22.716	ug/l	100
3) Chloromethane	1.300	50	44715	19.489	ug/l	99
4) Vinyl Chloride	1.380	62	48480	19.841	ug/l	99
5) Bromomethane	1.617	94	33887	22.441	ug/l	96
6) Chloroethane	1.697	64	30088	19.785	ug/l	96
7) Trichlorofluoromethane	1.898	101	71250	20.345	ug/l	99
8) Diethyl Ether	2.130	74	26662	19.074	ug/l	94
9) 1,1,2-Trichlorotrifluo...	2.337	101	39901	19.653	ug/l	97
10) Methyl Iodide	2.453	142	50587	17.502	ug/l	97
11) Tert butyl alcohol	2.928	59	39487	102.117	ug/l	99
12) 1,1-Dichloroethene	2.325	96	39681	19.150	ug/l	94
13) Acrolein	2.233	56	38865	127.548	ug/l	98
14) Allyl chloride	2.666	41	73696	21.184	ug/l	96
15) Acrylonitrile	3.050	53	121301	100.674	ug/l	98
16) Acetone	2.367	43	99898	108.954	ug/l	91
17) Carbon Disulfide	2.514	76	112473	18.691	ug/l	99
18) Methyl Acetate	2.690	43	55490	21.510	ug/l	96
19) Methyl tert-butyl Ether	3.099	73	137477	20.386	ug/l	100
20) Methylene Chloride	2.788	84	46445	20.266	ug/l	98
21) trans-1,2-Dichloroethene	3.087	96	38834	18.362	ug/l	97
22) Diisopropyl ether	3.745	45	138676	21.329	ug/l	99
23) Vinyl Acetate	3.709	43	614319	112.447	ug/l	95
24) 1,1-Dichloroethane	3.599	63	75325	20.775	ug/l	99
25) 2-Butanone	4.532	43	153848	113.689	ug/l #	89
26) 2,2-Dichloropropane	4.452	77	65032	20.271	ug/l	97
27) cis-1,2-Dichloroethene	4.471	96	47580	19.937	ug/l	94
28) Bromochloromethane	4.879	49	38829	21.849	ug/l	84
29) Tetrahydrofuran	4.971	42	103797	100.452	ug/l	95
30) Chloroform	5.068	83	78114	20.107	ug/l	93
31) Cyclohexane	5.452	56	60315	18.632	ug/l	99
32) 1,1,1-Trichloroethane	5.361	97	69881	20.075	ug/l	96
36) 1,1-Dichloropropene	5.672	75	50056	17.754	ug/l	99
37) Ethyl Acetate	4.690	43	72073	21.393	ug/l	96
38) Carbon Tetrachloride	5.659	117	61250	19.179	ug/l	93
39) Methylcyclohexane	7.360	83	57003	17.098	ug/l	97
40) Benzene	6.019	78	157310	18.821	ug/l	100

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120825\
 Data File : VX048745.D
 Acq On : 08 Dec 2025 11:44
 Operator : JC/MD
 Sample : VX1208WBS01
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 VX1208WBS01

Manual Integrations
 APPROVED

Reviewed By : John Carlone 12/09/2025
 Supervised By : Mahesh Dadoda 12/09/2025

Quant Time: Dec 09 04:06:52 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X120425W.M
 Quant Title : SW846 8260
 QLast Update : Fri Dec 05 03:27:34 2025
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Methacrylonitrile	4.891	41	31647	20.646	ug/l	94
42) 1,2-Dichloroethane	6.068	62	63162	21.351	ug/l	93
43) Isopropyl Acetate	6.312	43	105648	20.710	ug/l	94
44) Trichloroethene	7.104	130	38336	17.578	ug/l	94
45) 1,2-Dichloropropane	7.409	63	38148	18.261	ug/l	99
46) Dibromomethane	7.562	93	29948	18.904	ug/l	95
47) Bromodichloromethane	7.799	83	62155	18.818	ug/l	97
48) Methyl methacrylate	7.671	41	49411	19.255	ug/l	94
49) 1,4-Dioxane	7.641	88	11933	338.111	ug/l	95
51) 4-Methyl-2-Pentanone	8.549	43	333290	105.257	ug/l	96
52) Toluene	8.702	92	98537	19.551	ug/l	99
53) t-1,3-Dichloropropene	8.958	75	62916	20.022	ug/l	98
54) cis-1,3-Dichloropropene	8.348	75	65243	19.320	ug/l	94
55) 1,1,2-Trichloroethane	9.134	97	44019	21.707	ug/l	99
56) Ethyl methacrylate	9.098	69	67147	21.154	ug/l	97
57) 1,3-Dichloropropane	9.287	76	69273	20.543	ug/l	97
58) 2-Chloroethyl Vinyl ether	8.220	63	152119	88.668	ug/l	97
59) 2-Hexanone	9.409	43	232393	104.551	ug/l	98
60) Dibromochloromethane	9.500	129	53723	21.424	ug/l	100
61) 1,2-Dibromoethane	9.592	107	46114	21.010	ug/l	99
64) Tetrachloroethene	9.256	164	37431	16.498	ug/l	97
65) Chlorobenzene	10.061	112	118471	17.388	ug/l	98
66) 1,1,1,2-Tetrachloroethane	10.146	131	42596	17.479	ug/l	99
67) Ethyl Benzene	10.177	91	198185	17.826	ug/l	100
68) m/p-Xylenes	10.281	106	155889	35.995	ug/l	96
69) o-Xylene	10.622	106	73460	18.102	ug/l	97
70) Styrene	10.634	104	126903	18.215	ug/l	99
71) Bromoform	10.780	173	37658	17.309	ug/l #	99
73) Isopropylbenzene	10.945	105	188979	17.766	ug/l	99
74) N-amyl acetate	10.823	43	81515	17.637	ug/l	96
75) 1,1,2,2-Tetrachloroethane	11.195	83	63497	17.692	ug/l	99
76) 1,2,3-Trichloropropane	11.219	75	49341m	16.783	ug/l	
77) Bromobenzene	11.183	156	50590	17.819	ug/l	93
78) n-propylbenzene	11.286	91	218787	17.759	ug/l	98
79) 2-Chlorotoluene	11.347	91	131768	17.540	ug/l	97
80) 1,3,5-Trimethylbenzene	11.433	105	154012	17.615	ug/l	99
81) trans-1,4-Dichloro-2-b...	11.000	75	20596	16.641	ug/l	90
82) 4-Chlorotoluene	11.433	91	156772	17.883	ug/l	98
83) tert-Butylbenzene	11.695	119	158909	17.491	ug/l	98
84) 1,2,4-Trimethylbenzene	11.732	105	158196	17.937	ug/l	99
85) sec-Butylbenzene	11.872	105	190539	17.306	ug/l	98
86) p-Isopropyltoluene	11.988	119	164163	17.443	ug/l	100
87) 1,3-Dichlorobenzene	11.951	146	90505	17.292	ug/l	99
88) 1,4-Dichlorobenzene	12.024	146	92718	17.143	ug/l	99
89) n-Butylbenzene	12.311	91	144814	16.434	ug/l	99
90) Hexachloroethane	12.518	117	30759	16.013	ug/l	97
91) 1,2-Dichlorobenzene	12.317	146	85738	16.897	ug/l	98
92) 1,2-Dibromo-3-Chloropr...	12.920	75	13707	15.692	ug/l	95
93) 1,2,4-Trichlorobenzene	13.567	180	55758	16.638	ug/l	99
94) Hexachlorobutadiene	13.707	225	21763	15.700	ug/l	99
95) Naphthalene	13.756	128	177703	20.256	ug/l	100
96) 1,2,3-Trichlorobenzene	13.938	180	52737	17.064	ug/l	99

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120825\
Data File : VX048745.D
Acq On : 08 Dec 2025 11:44
Operator : JC/MD
Sample : VX1208WBS01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 5 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
VX1208WBS01

Manual Integrations
APPROVED

Reviewed By :John Carlone 12/09/2025
Supervised By :Mahesh Dadoda 12/09/2025

Quant Time: Dec 09 04:06:52 2025
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X120425W.M
Quant Title : SW846 8260
QLast Update : Fri Dec 05 03:27:34 2025
Response via : Initial Calibration

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

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

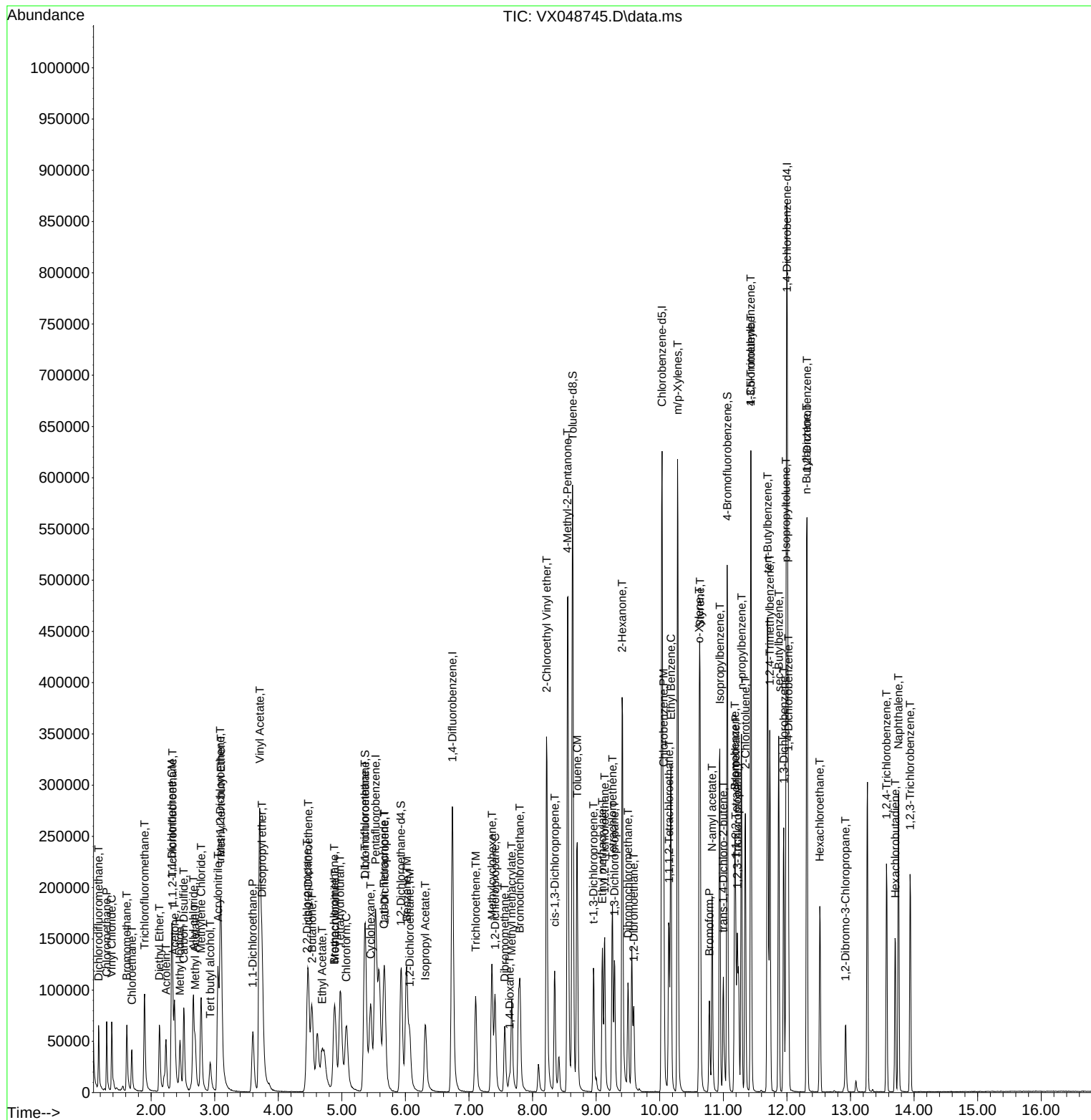
```
Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120825\
Data File : VX048745.D
Acq On    : 08 Dec 2025  11:44
Operator  : JC/MD
Sample    : VX1208WBS01
Misc      : 5.0mL/MSVOA_X/WATER
ALS Vial  : 5      Sample Multiplier: 1
```

Instrument :
MSVOA_X
ClientSampleId :
VX1208WBS01

Manual Integrations APPROVED

Reviewed By :John Carlone 12/09/2025
Supervised By :Mahesh Dadoda 12/09/2025

Quant Time: Dec 09 04:06:52 2025
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X120425W.M
Quant Title : SW846 8260
QLast Update : Fri Dec 05 03:27:34 2025
Response via : Initial Calibration





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

Report of Analysis

Client: JACOBS Engineering Group, Inc.
Project: Former Schlumberger STC PTC Site D3868221
Client Sample ID: VN1204WBSD01
Lab Sample ID: VN1204WBSD01
Analytical Method: 8260D
Sample Wt/Vol: 5 mL

Level: LOW
Final Vol: 5000 uL

Date Collected:
Date Received:
SDG No.: Q3757
Matrix: Water
% Solid: 0
Test: VOCMS Group3

CAS Number	Parameter	Conc.	Qua.	DF	MDL	LOQ / CRQL	Units	Date Ana.	BatchID
TARGETS									
75-01-4	Vinyl Chloride	19.0		1	0.26	1.00	ug/L	12/04/25 14:45	VN120425
75-35-4	1,1-Dichloroethene	19.5		1	0.23	1.00	ug/L	12/04/25 14:45	VN120425
75-34-3	1,1-Dichloroethane	19.7		1	0.23	1.00	ug/L	12/04/25 14:45	VN120425
156-59-2	cis-1,2-Dichloroethene	19.3		1	0.19	1.00	ug/L	12/04/25 14:45	VN120425
71-55-6	1,1,1-Trichloroethane	20.6		1	0.20	1.00	ug/L	12/04/25 14:45	VN120425
71-43-2	Benzene	19.3		1	0.15	1.00	ug/L	12/04/25 14:45	VN120425
107-06-2	1,2-Dichloroethane	21.6		1	0.22	1.00	ug/L	12/04/25 14:45	VN120425
79-01-6	Trichloroethene	18.7		1	0.090	1.00	ug/L	12/04/25 14:45	VN120425
79-00-5	1,1,2-Trichloroethane	20.7		1	0.21	1.00	ug/L	12/04/25 14:45	VN120425
127-18-4	Tetrachloroethene	17.3		1	0.23	1.00	ug/L	12/04/25 14:45	VN120425
SURROGATES									
17060-07-0	1,2-Dichloroethane-d4	56.0			70 (74) - 130 (125)	112%	SPK: 50		
1868-53-7	Dibromofluoromethane	53.1			70 (75) - 130 (124)	106%	SPK: 50		
2037-26-5	Toluene-d8	52.7			70 (86) - 130 (113)	105%	SPK: 50		
460-00-4	4-Bromofluorobenzene	56.3			70 (77) - 130 (121)	113%	SPK: 50		
INTERNAL STANDARDS									
		Area Count							
363-72-4	Pentafluorobenzene	291000							
540-36-3	1,4-Difluorobenzene	529000							
3114-55-4	Chlorobenzene-d5	470000							
3855-82-1	1,4-Dichlorobenzene-d4	230000							

U = Not Detected

LOQ = Limit of Quantitation

MDL = Method Detection Limit

LOD = Limit of Detection

E = Value Exceeds Calibration Range

Q = indicates LCS control criteria did not meet requirements

M = MS/MSD acceptance criteria did not meet requirements

J = Estimated Value

B = Analyte Found in Associated Method Blank

N = Presumptive Evidence of a Compound

* = Values outside of QC limits

D = Dilution

() = Laboratory InHouse Limit

A = Aldol-Condensation Reaction Products

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN120425\
 Data File : VN088444.D
 Acq On : 04 Dec 2025 14:45
 Operator : JC\MD
 Sample : VN1204WBSD01
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 VN1204WBSD01

Manual Integrations
 APPROVED

Reviewed By : John Carlone 12/05/2025
 Supervised By : Semsettin Yesilyurt 12/05/2025

Quant Time: Dec 05 00:15:47 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N112425W.M
 Quant Title : SW846 8260
 QLast Update : Tue Nov 25 01:34:52 2025
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.206	168	290727	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.077	114	529277	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.841	117	469866	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.770	152	229597	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.559	65	306737	56.007	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery = 112.020%			
35) Dibromofluoromethane	8.147	113	194821	53.116	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery = 106.240%			
50) Toluene-d8	10.541	98	719265	52.704	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery = 105.400%			
62) 4-Bromofluorobenzene	12.829	95	263766	56.328	ug/l	0.00
Spiked Amount 50.000	Range 77 - 121		Recovery = 112.660%			
Target Compounds					Qvalue	
2) Dichlorodifluoromethane	2.142	85	87020	19.928	ug/l	94
3) Chloromethane	2.383	50	84956	18.114	ug/l	97
4) Vinyl Chloride	2.536	62	90740	19.018	ug/l	96
5) Bromomethane	2.983	94	43017	18.733	ug/l	94
6) Chloroethane	3.148	64	55928	18.942	ug/l	99
7) Trichlorofluoromethane	3.518	101	138107	20.595	ug/l	99
8) Diethyl Ether	3.965	74	49843	19.803	ug/l	96
9) 1,1,2-Trichlorotrifluo...	4.377	101	71554	20.043	ug/l	99
10) Methyl Iodide	4.589	142	65248	18.463	ug/l	95
11) Tert butyl alcohol	5.518	59	80431	91.985	ug/l	99
12) 1,1-Dichloroethene	4.342	96	69022	19.460	ug/l	96
13) Acrolein	4.177	56	94599	113.005	ug/l	99
14) Allyl chloride	5.012	41	126430	19.231	ug/l	98
15) Acrylonitrile	5.706	53	250738	99.389	ug/l	98
16) Acetone	4.424	43	191583	96.727	ug/l	95
17) Carbon Disulfide	4.706	76	214025	18.466	ug/l	99
18) Methyl Acetate	5.018	43	126740	21.497	ug/l	99
19) Methyl tert-butyl Ether	5.783	73	277962	20.917	ug/l	100
20) Methylene Chloride	5.265	84	79618	18.788	ug/l	96
21) trans-1,2-Dichloroethene	5.771	96	74638	18.996	ug/l	84
22) Diisopropyl ether	6.659	45	291659	20.780	ug/l	97
23) Vinyl Acetate	6.589	43	1291270	114.189	ug/l	100
24) 1,1-Dichloroethane	6.553	63	156380	19.675	ug/l	99
25) 2-Butanone	7.465	43	318927	100.161	ug/l	97
26) 2,2-Dichloropropane	7.471	77	138156	20.137	ug/l	99
27) cis-1,2-Dichloroethene	7.471	96	88188	19.313	ug/l	96
28) Bromochloromethane	7.800	49	73205	18.489	ug/l	98
29) Tetrahydrofuran	7.824	42	234483	102.668	ug/l	98
30) Chloroform	7.947	83	157238	20.133	ug/l	93
31) Cyclohexane	8.241	56	140672	19.530	ug/l	97
32) 1,1,1-Trichloroethane	8.147	97	141532	20.610	ug/l	94
36) 1,1-Dichloropropene	8.353	75	108661	19.196	ug/l	99
37) Ethyl Acetate	7.547	43	150629	20.237	ug/l	99
38) Carbon Tetrachloride	8.341	117	114598	20.336	ug/l	98
39) Methylcyclohexane	9.577	83	114089	18.842	ug/l	97
40) Benzene	8.588	78	320551	19.294	ug/l	99

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN120425\
 Data File : VN088444.D
 Acq On : 04 Dec 2025 14:45
 Operator : JC\MD
 Sample : VN1204WBSD01
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 VN1204WBSD01

Manual Integrations
 APPROVED

Reviewed By :John Carlone 12/05/2025
 Supervised By :Semsettin Yesilyurt 12/05/2025

Quant Time: Dec 05 00:15:47 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N112425W.M
 Quant Title : SW846 8260
 QLast Update : Tue Nov 25 01:34:52 2025
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Methacrylonitrile	7.765	41	76908	20.862	ug/l	96
42) 1,2-Dichloroethane	8.653	62	136648	21.558	ug/l	99
43) Isopropyl Acetate	8.671	43	226089	19.881	ug/l	98
44) Trichloroethene	9.330	130	73312	18.675	ug/l	89
45) 1,2-Dichloropropane	9.600	63	83457	19.606	ug/l	90
46) Dibromomethane	9.688	93	61573	20.269	ug/l	98
47) Bromodichloromethane	9.865	83	126000	20.328	ug/l	98
48) Methyl methacrylate	9.659	41	106808	20.584	ug/l	98
49) 1,4-Dioxane	9.682	88	23459	355.507	ug/l #	99
51) 4-Methyl-2-Pentanone	10.424	43	707725	105.581	ug/l	99
52) Toluene	10.606	92	192863	20.104	ug/l	99
53) t-1,3-Dichloropropene	10.812	75	127987	20.116	ug/l	93
54) cis-1,3-Dichloropropene	10.288	75	132808	19.849	ug/l	98
55) 1,1,2-Trichloroethane	10.994	97	76705	20.666	ug/l	98
56) Ethyl methacrylate	10.859	69	126407	21.560	ug/l	99
57) 1,3-Dichloropropane	11.141	76	138389	20.668	ug/l	100
58) 2-Chloroethyl Vinyl ether	10.141	63	320414	94.794	ug/l	98
59) 2-Hexanone	11.177	43	425614	100.793	ug/l	96
60) Dibromochloromethane	11.335	129	88166	20.837	ug/l	99
61) 1,2-Dibromoethane	11.447	107	78440	20.870	ug/l	99
64) Tetrachloroethene	11.082	164	64323	17.268	ug/l	97
65) Chlorobenzene	11.871	112	208506	19.756	ug/l	100
66) 1,1,1,2-Tetrachloroethane	11.935	131	71468	20.077	ug/l	99
67) Ethyl Benzene	11.941	91	364543	19.762	ug/l	98
68) m/p-Xylenes	12.047	106	273590	40.003	ug/l	97
69) o-Xylene	12.376	106	129944	19.929	ug/l	100
70) Styrene	12.388	104	216021	20.404	ug/l	98
71) Bromoform	12.559	173	53817	20.641	ug/l #	99
73) Isopropylbenzene	12.671	105	347231	20.454	ug/l	98
74) N-amyl acetate	12.682	43	111327m	19.256	ug/l	
75) 1,1,2,2-Tetrachloroethane	12.918	83	111054	20.235	ug/l	99
76) 1,2,3-Trichloropropane	12.971	75	113801m	21.162	ug/l	
77) Bromobenzene	12.959	156	77652	19.992	ug/l	96
78) n-propylbenzene	13.012	91	421199	20.281	ug/l	100
79) 2-Chlorotoluene	13.106	91	253171	19.962	ug/l	100
80) 1,3,5-Trimethylbenzene	13.153	105	281269	19.995	ug/l	99
81) trans-1,4-Dichloro-2-b...	12.718	75	35239	18.590	ug/l #	93
82) 4-Chlorotoluene	13.200	91	267452	21.232	ug/l	99
83) tert-Butylbenzene	13.412	119	239926	20.064	ug/l	98
84) 1,2,4-Trimethylbenzene	13.459	105	286265	20.474	ug/l	97
85) sec-Butylbenzene	13.594	105	344400	20.048	ug/l	99
86) p-Isopropyltoluene	13.706	119	283564	20.130	ug/l	98
87) 1,3-Dichlorobenzene	13.712	146	148014	19.576	ug/l	99
88) 1,4-Dichlorobenzene	13.788	146	152072	19.200	ug/l	96
89) n-Butylbenzene	14.035	91	267202	19.536	ug/l	97
90) Hexachloroethane	14.312	117	48757	18.947	ug/l	99
91) 1,2-Dichlorobenzene	14.082	146	144112	20.254	ug/l	99
92) 1,2-Dibromo-3-Chloropr...	14.694	75	26992	20.271	ug/l	98
93) 1,2,4-Trichlorobenzene	15.364	180	81616	19.305	ug/l	97
94) Hexachlorobutadiene	15.470	225	27309	17.976	ug/l	94
95) Naphthalene	15.612	128	302681	20.661	ug/l	99
96) 1,2,3-Trichlorobenzene	15.812	180	77952	18.946	ug/l	96

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN120425\
Data File : VN088444.D
Acq On : 04 Dec 2025 14:45
Operator : JC\MD
Sample : VN1204WBSD01
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
VN1204WBSD01

Manual Integrations
APPROVED

Reviewed By :John Carlone 12/05/2025
Supervised By :Semsettin Yesilyurt 12/05/2025

Quant Time: Dec 05 00:15:47 2025
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N112425W.M
Quant Title : SW846 8260
QLast Update : Tue Nov 25 01:34:52 2025
Response via : Initial Calibration

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

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

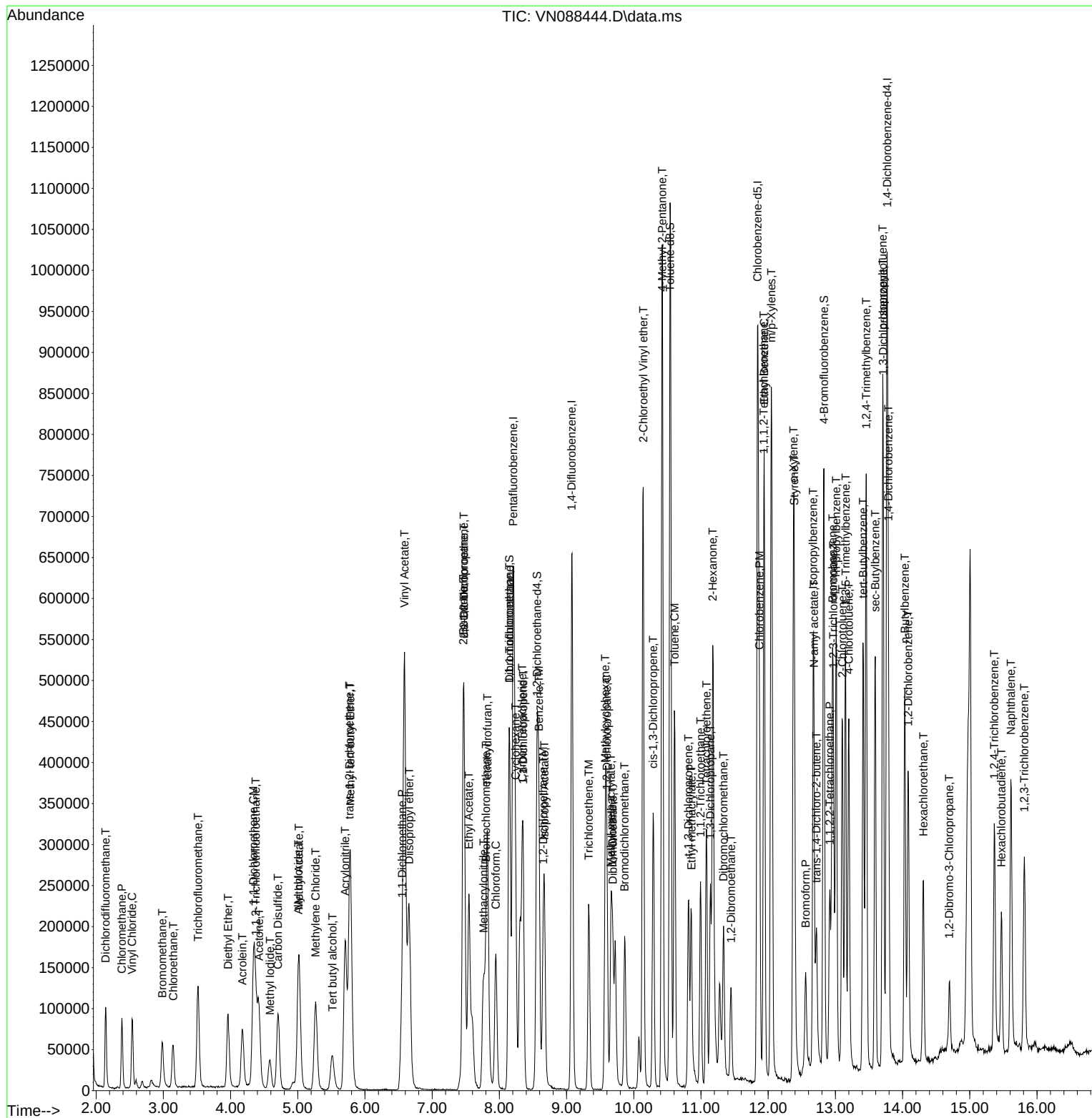
```
Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN120425\
Data File : VN088444.D
Acq On    : 04 Dec 2025  14:45
Operator  : JC\MD
Sample    : VN1204WBSD01
Misc      : 5.0mL/MSVOA_N/WATER
ALS Vial  : 10   Sample Multiplier: 1
```

Instrument :
MSVOA_N
ClientSampleId :
VN1204WBSD01

Manual Integrations APPROVED

Reviewed By :John Carlone 12/05/2025
Supervised By :Semsettin Yesilyurt 12/05/2025

Quant Time: Dec 05 00:15:47 2025
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N112425W.M
Quant Title : SW846 8260
QLast Update : Tue Nov 25 01:34:52 2025
Response via : Initial Calibration



Manual Integration Report

Sequence:	VN112425	Instrument	MSVOA_n
-----------	----------	------------	---------

Sample ID	File ID	Parameter	Review By	Review On	Supervised By	Supervised On	Reason
VSTDICC001	VN088344.D	1,1-Dichloroethene	JOHN	11/25/2025 9:35:55 AM	MMDadoda	11/26/2025 1:51:48 PM	Peak Integrated by Software
VSTDICC001	VN088344.D	1,2,3-Trichloropropane	JOHN	11/25/2025 9:35:55 AM	MMDadoda	11/26/2025 1:51:48 PM	Peak Integrated by Software
VSTDICC001	VN088344.D	1,4-Dichlorobenzene	JOHN	11/25/2025 9:35:55 AM	MMDadoda	11/26/2025 1:51:48 PM	Peak Integrated by Software
VSTDICC001	VN088344.D	2-Hexanone	JOHN	11/25/2025 9:35:55 AM	MMDadoda	11/26/2025 1:51:48 PM	Peak Integrated by Software
VSTDICC001	VN088344.D	cis-1,2-Dichloroethene	JOHN	11/25/2025 9:35:55 AM	MMDadoda	11/26/2025 1:51:48 PM	Peak Integrated by Software
VSTDICC001	VN088344.D	Diisopropyl ether	JOHN	11/25/2025 9:35:55 AM	MMDadoda	11/26/2025 1:51:48 PM	Peak Integrated by Software
VSTDICC005	VN088345.D	1,2,3-Trichloropropane	JOHN	11/25/2025 9:36:00 AM	MMDadoda	11/26/2025 1:51:51 PM	Peak Integrated by Software
VSTDICC005	VN088345.D	Vinyl Acetate	JOHN	11/25/2025 9:36:00 AM	MMDadoda	11/26/2025 1:51:51 PM	Peak Integrated by Software
VSTDICC020	VN088346.D	1,2,3-Trichloropropane	JOHN	11/25/2025 9:36:04 AM	MMDadoda	11/26/2025 1:51:55 PM	Peak Integrated by Software
VSTDICC020	VN088346.D	N-amyl acetate	JOHN	11/25/2025 9:36:04 AM	MMDadoda	11/26/2025 1:51:55 PM	Peak Integrated by Software
VSTDICC020	VN088346.D	Vinyl Acetate	JOHN	11/25/2025 9:36:04 AM	MMDadoda	11/26/2025 1:51:55 PM	Peak Integrated by Software
VSTDICCC050	VN088347.D	1,2,3-Trichloropropane	JOHN	11/25/2025 9:36:10 AM	MMDadoda	11/26/2025 1:51:58 PM	Peak Integrated by Software
VSTDICCC050	VN088347.D	Vinyl Acetate	JOHN	11/25/2025 9:36:10 AM	MMDadoda	11/26/2025 1:51:58 PM	Peak Integrated by Software

Manual Integration Report

Sequence:

VN112425

Instrument

MSVOA_n

Sample ID	File ID	Parameter	Review By	Review On	Supervised By	Supervised On	Reason
VSTDICC100	VN088348.D	1,2,3-Trichloropropane	JOHN	11/25/2025 9:36:15 AM	MMDadoda	11/26/2025 1:52:01 PM	Peak Integrated by Software
VSTDICC100	VN088348.D	Allyl chloride	JOHN	11/25/2025 9:36:15 AM	MMDadoda	11/26/2025 1:52:01 PM	Peak Integrated by Software
VSTDICC100	VN088348.D	Vinyl Acetate	JOHN	11/25/2025 9:36:15 AM	MMDadoda	11/26/2025 1:52:01 PM	Peak Integrated by Software
VSTDICC150	VN088349.D	1,2,3-Trichloropropane	JOHN	11/25/2025 9:36:20 AM	MMDadoda	11/26/2025 1:52:04 PM	Peak Integrated by Software
VSTDICC150	VN088349.D	Allyl chloride	JOHN	11/25/2025 9:36:20 AM	MMDadoda	11/26/2025 1:52:04 PM	Peak Integrated by Software
VSTDICC150	VN088349.D	Vinyl Acetate	JOHN	11/25/2025 9:36:20 AM	MMDadoda	11/26/2025 1:52:04 PM	Peak Integrated by Software
VSTDICV050	VN088351.D	1,2,3-Trichloropropane	JOHN	11/25/2025 9:36:25 AM	MMDadoda	11/26/2025 1:52:07 PM	Peak Integrated by Software
VSTDICV050	VN088351.D	N-amyl acetate	JOHN	11/25/2025 9:36:25 AM	MMDadoda	11/26/2025 1:52:07 PM	Peak Integrated by Software
VSTDICV050	VN088351.D	Vinyl Acetate	JOHN	11/25/2025 9:36:25 AM	MMDadoda	11/26/2025 1:52:07 PM	Peak Integrated by Software

Manual Integration Report

Sequence:	VN120425	Instrument	MSVOA_n
-----------	----------	------------	---------

Sample ID	File ID	Parameter	Review By	Review On	Supervised By	Supervised On	Reason
VSTDCCC050	VN088437.D	1,2,3-Trichloropropane	JOHN	12/5/2025 8:18:14 AM	sam	12/5/2025 2:18:25 PM	Peak Integrated by Software
VSTDCCC050	VN088437.D	N-amyl acetate	JOHN	12/5/2025 8:18:14 AM	sam	12/5/2025 2:18:25 PM	Peak Integrated by Software
VN1204WBS01	VN088443.D	1,2,3-Trichloropropane	JOHN	12/5/2025 8:18:23 AM	sam	12/5/2025 2:18:38 PM	Peak Integrated by Software
VN1204WBS01	VN088443.D	N-amyl acetate	JOHN	12/5/2025 8:18:23 AM	sam	12/5/2025 2:18:38 PM	Peak Integrated by Software
VN1204WBSD01	VN088444.D	1,2,3-Trichloropropane	JOHN	12/5/2025 8:18:28 AM	sam	12/5/2025 2:18:43 PM	Peak Integrated by Software
VN1204WBSD01	VN088444.D	N-amyl acetate	JOHN	12/5/2025 8:18:28 AM	sam	12/5/2025 2:18:43 PM	Peak Integrated by Software
Q3757-11	VN088450.D	Acetone	JOHN	12/5/2025 8:18:38 AM	sam	12/5/2025 2:19:00 PM	Peak Integrated by Software
VSTDCCC050	VN088461.D	1,2,3-Trichloropropane	JOHN	12/5/2025 8:18:43 AM	sam	12/5/2025 2:18:57 PM	Peak Integrated by Software
VSTDCCC050	VN088461.D	N-amyl acetate	JOHN	12/5/2025 8:18:43 AM	sam	12/5/2025 2:18:57 PM	Peak Integrated by Software

Manual Integration Report

Sequence:	VX120425	Instrument	MSVOA_x
-----------	----------	------------	---------

Sample ID	File ID	Parameter	Review By	Review On	Supervised By	Supervised On	Reason
VSTDICC001	VX048714.D	1,2,3-Trichloropropane	JOHN	12/5/2025 8:04:03 AM	sam	12/5/2025 2:19:32 PM	Peak Integrated by Software
VSTDICC001	VX048714.D	1,4-Dichlorobenzene	JOHN	12/5/2025 8:04:03 AM	sam	12/5/2025 2:19:32 PM	Peak Integrated by Software
VSTDICC001	VX048714.D	Ethyl Acetate	JOHN	12/5/2025 8:04:03 AM	sam	12/5/2025 2:19:32 PM	Peak Integrated by Software
VSTDICC005	VX048715.D	1,2,3-Trichloropropane	JOHN	12/5/2025 8:04:10 AM	sam	12/5/2025 2:19:37 PM	Peak Integrated by Software
VSTDICC005	VX048715.D	1,2-Dichloroethane-d4	JOHN	12/5/2025 8:04:10 AM	sam	12/5/2025 2:19:37 PM	Peak Integrated by Software
VSTDICC005	VX048715.D	4-Bromofluorobenzene	JOHN	12/5/2025 8:04:10 AM	sam	12/5/2025 2:19:37 PM	Peak Integrated by Software
VSTDICC005	VX048715.D	Acrolein	JOHN	12/5/2025 8:04:10 AM	sam	12/5/2025 2:19:37 PM	Peak Integrated by Software
VSTDICC005	VX048715.D	Bromomethane	JOHN	12/5/2025 8:04:10 AM	sam	12/5/2025 2:19:37 PM	Peak Integrated by Software
VSTDICC005	VX048715.D	Cyclohexane	JOHN	12/5/2025 8:04:10 AM	sam	12/5/2025 2:19:37 PM	Peak Integrated by Software
VSTDICC005	VX048715.D	Dibromofluoromethane	JOHN	12/5/2025 8:04:10 AM	sam	12/5/2025 2:19:37 PM	Peak Integrated by Software
VSTDICC005	VX048715.D	Ethyl Acetate	JOHN	12/5/2025 8:04:10 AM	sam	12/5/2025 2:19:37 PM	Peak Integrated by Software
VSTDICC005	VX048715.D	Methyl Acetate	JOHN	12/5/2025 8:04:10 AM	sam	12/5/2025 2:19:37 PM	Peak Integrated by Software
VSTDICC005	VX048715.D	Methyl Iodide	JOHN	12/5/2025 8:04:10 AM	sam	12/5/2025 2:19:37 PM	Peak Integrated by Software

Manual Integration Report

Sequence:

VX120425

Instrument

MSVOA_x

Sample ID	File ID	Parameter	Review By	Review On	Supervised By	Supervised On	Reason
VSTDIC005	VX048715.D	Tert butyl alcohol	JOHN	12/5/2025 8:04:10 AM	sam	12/5/2025 2:19:37 PM	Peak Integrated by Software
VSTDIC005	VX048715.D	Toluene-d8	JOHN	12/5/2025 8:04:10 AM	sam	12/5/2025 2:19:37 PM	Peak Integrated by Software
VSTDIC005	VX048715.D	trans-1,4-Dichloro-2-but ene	JOHN	12/5/2025 8:04:10 AM	sam	12/5/2025 2:19:37 PM	Peak Integrated by Software
VSTDIC020	VX048716.D	1,2,3-Trichloropropane	JOHN	12/5/2025 8:04:16 AM	sam	12/5/2025 2:19:42 PM	Peak Integrated by Software
VSTDIC050	VX048717.D	1,2,3-Trichloropropane	JOHN	12/5/2025 8:04:22 AM	sam	12/5/2025 2:20:03 PM	Peak Integrated by Software
VSTDIC100	VX048718.D	1,2,3-Trichloropropane	JOHN	12/5/2025 8:04:27 AM	sam	12/5/2025 2:19:47 PM	Peak Integrated by Software
VSTDIC150	VX048719.D	1,2,3-Trichloropropane	JOHN	12/5/2025 8:04:32 AM	sam	12/5/2025 2:20:07 PM	Peak Integrated by Software
VSTDICV050	VX048721.D	1,2,3-Trichloropropane	JOHN	12/5/2025 8:04:41 AM	sam	12/5/2025 2:19:52 PM	Peak Integrated by Software

Manual Integration Report

Sequence:	VX120525	Instrument	MSVOA_x
-----------	----------	------------	---------

Sample ID	File ID	Parameter	Review By	Review On	Supervised By	Supervised On	Reason
VSTDCCC050	VX048723.D	1,2,3-Trichloropropane	JOHN	12/8/2025 8:07:20 AM	sam	12/8/2025 11:02:20 AM	Peak Integrated by Software
VX1205WBS01	VX048726.D	1,2,3-Trichloropropane	JOHN	12/8/2025 8:07:25 AM	sam	12/8/2025 11:02:26 AM	Peak Integrated by Software
VSTDCCC050	VX048740.D	1,2,3-Trichloropropane	JOHN	12/8/2025 8:07:34 AM	sam	12/8/2025 11:02:36 AM	Peak Integrated by Software

Manual Integration Report

Sequence:	VX120825	Instrument	MSVOA_x
-----------	----------	------------	---------

Sample ID	File ID	Parameter	Review By	Review On	Supervised By	Supervised On	Reason
VSTDCCC050	VX048742.D	1,2,3-Trichloropropane	JOHN	12/9/2025 8:06:28 AM	MMDadoda	12/9/2025 12:23:37 PM	Peak Integrated by Software
VX1208WBS01	VX048745.D	1,2,3-Trichloropropane	JOHN	12/9/2025 8:06:33 AM	MMDadoda	12/9/2025 12:23:40 PM	Peak Integrated by Software
VSTDCCC050	VX048766.D	1,2,3-Trichloropropane	JOHN	12/9/2025 8:06:52 AM	MMDadoda	12/9/2025 12:23:58 PM	Peak Integrated by Software

Instrument ID: **MSVOA_N**

Daily Analysis Runlog For Sequence/QC Batch ID # VN112425

Review By	John Carlone	Review On	11/25/2025 9:36:43 AM		
Supervise By	Maresh Dadoda	Supervise On	11/26/2025 1:52:19 PM		
SubDirectory	VN112425	HP Acquire Method	MSVOA_N	HP Processing Method	82N112425W.M
STD. NAME		STD REF.#			
Tune/Reschk		VP136378			
Initial Calibration Stds		VP136444,VP136445,VP136446,VP136447,VP136448,VP136449			
CCC					
Internal Standard/PEM					
ICV/I.BLK		VP136450			
Surrogate Standard					
MS/MSD Standard					
LCS Standard					

Sr#	SampleId	Data File Name	Date-Time	Operator	Status
1	BFB	VN088343.D	24 Nov 2025 08:17	JC\MD	Ok
2	VSTDIC001	VN088344.D	24 Nov 2025 12:38	JC\MD	Ok,M
3	VSTDIC005	VN088345.D	24 Nov 2025 13:11	JC\MD	Ok,M
4	VSTDIC020	VN088346.D	24 Nov 2025 13:32	JC\MD	Ok,M
5	VSTDIC050	VN088347.D	24 Nov 2025 13:53	JC\MD	Ok,M
6	VSTDIC100	VN088348.D	24 Nov 2025 14:14	JC\MD	Ok,M
7	VSTDIC150	VN088349.D	24 Nov 2025 14:35	JC\MD	Ok,M
8	IBLK	VN088350.D	24 Nov 2025 14:56	JC\MD	Ok
9	VSTDICV050	VN088351.D	24 Nov 2025 15:18	JC\MD	Ok,M

M : Manual Integration

Instrument ID: **MSVOA_N**

Daily Analysis Runlog For Sequence/QC Batch ID # VN120425

Review By	John Carlone	Review On	12/5/2025 8:23:32 AM
Supervise By	Semsettin Yesilyurt	Supervise On	12/5/2025 2:14:29 PM
SubDirectory	VN120425	HP Acquire Method	HP Processing Method 82N112425W.M
STD. NAME	STD REF.#		
Tune/Reschk Initial Calibration Stds	VP136505		
CCC Internal Standard/PEM ICV/I.BLK Surrogate Standard MS/MSD Standard LCS Standard	VP136506,VP136507		

Sr#	SampleId	Data File Name	Date-Time	Operator	Status
1	BFB	VN088436.D	04 Dec 2025 11:03	JC\MD	Ok
2	VSTDCCC050	VN088437.D	04 Dec 2025 11:44	JC\MD	Ok,M
3	VN1204MBL01	VN088438.D	04 Dec 2025 12:17	JC\MD	Ok
4	VN1204WBL01	VN088439.D	04 Dec 2025 12:38	JC\MD	Ok
5	VN1204MBS01	VN088440.D	04 Dec 2025 13:11	JC\MD	Ok,M
6	Q3768-01	VN088441.D	04 Dec 2025 13:31	JC\MD	Not Ok
7	IBLK	VN088442.D	04 Dec 2025 13:52	JC\MD	Ok
8	VN1204WBS01	VN088443.D	04 Dec 2025 14:12	JC\MD	Ok,M
9	VN1204WBSD01	VN088444.D	04 Dec 2025 14:45	JC\MD	Ok,M
10	Q3756-03	VN088445.D	04 Dec 2025 15:06	JC\MD	Ok
11	Q3756-01	VN088446.D	04 Dec 2025 15:26	JC\MD	Ok
12	Q3756-05	VN088447.D	04 Dec 2025 15:47	JC\MD	Ok
13	Q3756-07	VN088448.D	04 Dec 2025 16:07	JC\MD	Ok
14	Q3757-12	VN088449.D	04 Dec 2025 16:28	JC\MD	Ok
15	Q3757-11	VN088450.D	04 Dec 2025 16:48	JC\MD	Ok,M
16	Q3757-01	VN088451.D	04 Dec 2025 17:09	JC\MD	Not Ok
17	Q3757-03	VN088452.D	04 Dec 2025 17:29	JC\MD	Not Ok
18	Q3757-05	VN088453.D	04 Dec 2025 17:50	JC\MD	Not Ok
19	Q3757-07	VN088454.D	04 Dec 2025 18:10	JC\MD	Not Ok
20	Q3757-09	VN088455.D	04 Dec 2025 18:31	JC\MD	Not Ok
21	Q3750-07	VN088456.D	04 Dec 2025 18:51	JC\MD	Ok

Instrument ID: **MSVOA_N**

Daily Analysis Runlog For Sequence/QC Batch ID # VN120425

Review By	John Carlone	Review On	12/5/2025 8:23:32 AM
Supervise By	Semsettin Yesilyurt	Supervise On	12/5/2025 2:14:29 PM
SubDirectory	VN120425	HP Acquire Method	HP Processing Method 82N112425W.M
STD. NAME	STD REF.#		
Tune/Reschk Initial Calibration Stds	VP136505		
CCC Internal Standard/PEM ICV/I.BLK Surrogate Standard MS/MSD Standard LCS Standard	VP136506,VP136507		

22	Q3750-08	VN088457.D	04 Dec 2025 19:11	JC\MD	Ok
23	PB170809TB	VN088458.D	04 Dec 2025 19:32	JC\MD	Ok
24	Q3775-01	VN088459.D	04 Dec 2025 19:52	JC\MD	Not Ok
25	IBLK	VN088460.D	04 Dec 2025 20:13	JC\MD	Ok
26	VSTDCCC050	VN088461.D	04 Dec 2025 20:33	JC\MD	Ok,M

M : Manual Integration

Instrument ID: **MSVOA_X**

Daily Analysis Runlog For Sequence/QC Batch ID # VX120425

Review By	John Carlone	Review On	12/5/2025 8:05:59 AM
Supervise By	Semsettin Yesilyurt	Supervise On	12/5/2025 2:14:46 PM
SubDirectory	VX120425	HP Acquire Method	HP Processing Method 82X120425
STD. NAME	STD REF.#		
Tune/Reschk	VP136497		
Initial Calibration Stds	VP136526,VP136527,VP136528,VP136529,VP136530,VP136531		
CCC			
Internal Standard/PEM			
ICV/I.BLK	VP136532		
Surrogate Standard			
MS/MSD Standard			
LCS Standard			

Sr#	SampleId	Data File Name	Date-Time	Operator	Status
1	BFB	VX048713.D	04 Dec 2025 08:15	JC/MD	Ok
2	VSTDIC001	VX048714.D	04 Dec 2025 08:47	JC/MD	Ok,M
3	VSTDIC005	VX048715.D	04 Dec 2025 09:16	JC/MD	Ok,M
4	VSTDIC020	VX048716.D	04 Dec 2025 09:36	JC/MD	Ok,M
5	VSTDIC050	VX048717.D	04 Dec 2025 09:57	JC/MD	Ok,M
6	VSTDIC100	VX048718.D	04 Dec 2025 10:17	JC/MD	Ok,M
7	VSTDIC150	VX048719.D	04 Dec 2025 10:38	JC/MD	Ok,M
8	IBLK	VX048720.D	04 Dec 2025 10:58	JC/MD	Ok
9	VSTDICV050	VX048721.D	04 Dec 2025 14:04	JC/MD	Ok,M

M : Manual Integration

Instrument ID: **MSVOA_X**

Daily Analysis Runlog For Sequence/QC Batch ID # VX120525

Review By	John Carlone	Review On	12/8/2025 8:24:07 AM
Supervise By	Semsettin Yesilyurt	Supervise On	12/8/2025 11:02:48 AM
SubDirectory	VX120525	HP Acquire Method	HP Processing Method 82X120425
STD. NAME	STD REF.#		
Tune/Reschk Initial Calibration Stds	VP136523		
CCC Internal Standard/PEM ICV/I.BLK Surrogate Standard MS/MSD Standard LCS Standard	VP136524,VP136525		

Sr#	SampleId	Data File Name	Date-Time	Operator	Status
1	BFB	VX048722.D	05 Dec 2025 08:14	JC/MD	Ok
2	VSTDCCC050	VX048723.D	05 Dec 2025 08:45	JC/MD	Ok,M
3	VX1205MBL01	VX048724.D	05 Dec 2025 09:14	JC/MD	Ok
4	VX1205WBL01	VX048725.D	05 Dec 2025 10:34	JC/MD	Ok
5	VX1205WBS01	VX048726.D	05 Dec 2025 11:31	JC/MD	Ok,M
6	VX1205WBSD01	VX048727.D	05 Dec 2025 11:59	JC/MD	Ok,M
7	Q3775-01	VX048728.D	05 Dec 2025 12:20	JC/MD	ReRun
8	Q3775-01RE	VX048729.D	05 Dec 2025 12:50	JC/MD	Confirms
9	IBLK	VX048730.D	05 Dec 2025 13:10	JC/MD	Ok
10	Q3757-01	VX048731.D	05 Dec 2025 13:32	JC/MD	Ok
11	Q3757-03	VX048732.D	05 Dec 2025 13:53	JC/MD	Ok
12	Q3757-05	VX048733.D	05 Dec 2025 14:13	JC/MD	ReRun
13	Q3757-07	VX048734.D	05 Dec 2025 14:33	JC/MD	Ok
14	Q3757-09	VX048735.D	05 Dec 2025 14:54	JC/MD	ReRun
15	Q3789-01	VX048736.D	05 Dec 2025 15:14	JC/MD	Ok
16	Q3789-02	VX048737.D	05 Dec 2025 15:35	JC/MD	Ok
17	Q3789-03	VX048738.D	05 Dec 2025 15:55	JC/MD	ReRun
18	Q3789-04	VX048739.D	05 Dec 2025 16:15	JC/MD	ReRun
19	VSTDCCC050	VX048740.D	05 Dec 2025 16:36	JC/MD	Ok,M

M : Manual Integration

Instrument ID: **MSVOA_X**

Daily Analysis Runlog For Sequence/QC Batch ID # VX120825

Review By	John Carlone	Review On	12/9/2025 8:11:51 AM
Supervise By	Maresh Dadoda	Supervise On	12/9/2025 12:45:16 PM
SubDirectory	VX120825	HP Acquire Method	HP Processing Method 82X120425W.M
STD. NAME	STD REF.#		
Tune/Reschk Initial Calibration Stds	VP136561		
CCC Internal Standard/PEM ICV/I.BLK Surrogate Standard MS/MSD Standard LCS Standard	VP136570,VP136571		

Sr#	SampleId	Data File Name	Date-Time	Operator	Status
1	BFB	VX048741.D	08 Dec 2025 08:17	JC/MD	Ok
2	VSTDCCC050	VX048742.D	08 Dec 2025 09:20	JC/MD	Ok,M
3	VX1208MBL01	VX048743.D	08 Dec 2025 09:49	JC/MD	Ok
4	VX1208WBL01	VX048744.D	08 Dec 2025 11:15	JC/MD	Ok
5	VX1208WBS01	VX048745.D	08 Dec 2025 11:44	JC/MD	Ok,M
6	VX1208WBSD01	VX048746.D	08 Dec 2025 12:12	JC/MD	Ok,M
7	Q3757-05RE	VX048747.D	08 Dec 2025 12:33	JC/MD	Confirms
8	Q3757-09RE	VX048748.D	08 Dec 2025 12:53	JC/MD	Confirms
9	Q3787-01	VX048749.D	08 Dec 2025 13:14	JC/MD	Ok
10	Q3787-02MS	VX048750.D	08 Dec 2025 13:35	JC/MD	Not Ok
11	Q3787-03MSD	VX048751.D	08 Dec 2025 13:55	JC/MD	Not Ok
12	Q3787-07	VX048752.D	08 Dec 2025 14:16	JC/MD	ReRun
13	Q3787-09	VX048753.D	08 Dec 2025 14:36	JC/MD	ReRun
14	Q3787-11	VX048754.D	08 Dec 2025 14:57	JC/MD	ReRun
15	Q3787-13	VX048755.D	08 Dec 2025 15:18	JC/MD	ReRun
16	Q3787-15	VX048756.D	08 Dec 2025 15:38	JC/MD	Ok
17	Q3787-17	VX048757.D	08 Dec 2025 15:59	JC/MD	ReRun
18	Q3787-19	VX048758.D	08 Dec 2025 16:19	JC/MD	ReRun
19	Q3787-21	VX048759.D	08 Dec 2025 16:40	JC/MD	ReRun
20	Q3788-07	VX048760.D	08 Dec 2025 17:00	JC/MD	ReRun
21	Q3788-09	VX048761.D	08 Dec 2025 17:21	JC/MD	ReRun

Instrument ID: **MSVOA_X**

Daily Analysis Runlog For Sequence/QC Batch ID # VX120825

Review By	John Carlone	Review On	12/9/2025 8:11:51 AM
Supervise By	Mahesh Dadoda	Supervise On	12/9/2025 12:45:16 PM
SubDirectory	VX120825	HP Acquire Method	HP Processing Method 82X120425W.M
STD. NAME	STD REF.#		
Tune/Reschk Initial Calibration Stds	VP136561		
CCC Internal Standard/PEM ICV/I.BLK Surrogate Standard MS/MSD Standard LCS Standard	VP136570,VP136571		

22	Q3788-11	VX048762.D	08 Dec 2025 17:41	JC/MD	ReRun
23	Q3788-13	VX048763.D	08 Dec 2025 18:02	JC/MD	Ok
24	Q3787-02MS	VX048764.D	08 Dec 2025 18:22	JC/MD	Ok,M
25	Q3787-03MSD	VX048765.D	08 Dec 2025 18:43	JC/MD	Ok,M
26	VSTDCCC050	VX048766.D	08 Dec 2025 19:03	JC/MD	Ok,M

M : Manual Integration

Instrument ID: MSVOA_N

Daily Analysis Runlog For Sequence/QC Batch ID # VN112425

Review By	John Carlone	Review On	11/25/2025 9:36:43 AM
Supervise By	Maresh Dadoda	Supervise On	11/26/2025 1:52:19 PM
SubDirectory	VN112425	HP Acquire Method	MSVOA_N HP Processing Method 82N112425W.M

STD. NAME	STD REF.#
Tune/Reschk	VP136378
Initial Calibration Stds	VP136444,VP136445,VP136446,VP136447,VP136448,VP136449
CCC	
Internal Standard/PEM	
ICV/I.BLK	VP136450
Surrogate Standard	
MS/MSD Standard	
LCS Standard	

Sr#	SampleID	ClientID	Data File Name	Date-Time	Comment	Operator	Status
1	BFB	BFB	VN088343.D	24 Nov 2025 08:17		JC\MD	Ok
2	VSTDICC001	VSTDICC001	VN088344.D	24 Nov 2025 12:38		JC\MD	Ok,M
3	VSTDICC005	VSTDICC005	VN088345.D	24 Nov 2025 13:11	COM.#74 Peak not detected in 1 and 5 PPB	JC\MD	Ok,M
4	VSTDICC020	VSTDICC020	VN088346.D	24 Nov 2025 13:32		JC\MD	Ok,M
5	VSTDICCC050	VSTDICCC050	VN088347.D	24 Nov 2025 13:53	Method failed for Comp. #74	JC\MD	Ok,M
6	VSTDICC100	VSTDICC100	VN088348.D	24 Nov 2025 14:14		JC\MD	Ok,M
7	VSTDICC150	VSTDICC150	VN088349.D	24 Nov 2025 14:35		JC\MD	Ok,M
8	IBLK	IBLK	VN088350.D	24 Nov 2025 14:56		JC\MD	Ok
9	VSTDICV050	ICVVN112425	VN088351.D	24 Nov 2025 15:18		JC\MD	Ok,M

M : Manual Integration

Instrument ID: MSVOA_N

Daily Analysis Runlog For Sequence/QC Batch ID # VN120425

Review By	John Carlone	Review On	12/5/2025 8:23:32 AM
Supervise By	Semsettin Yesilyurt	Supervise On	12/5/2025 2:14:29 PM
SubDirectory	VN120425	HP Acquire Method	HP Processing Method 82N112425W.M

STD. NAME	STD REF.#
Tune/Reschk Initial Calibration Stds	VP136505
CCC Internal Standard/PEM ICV/I.BLK Surrogate Standard MS/MSD Standard LCS Standard	VP136506,VP136507

Sr#	SampleID	ClientID	Data File Name	Date-Time	Comment	Operator	Status
1	BFB	BFB	VN088436.D	04 Dec 2025 11:03		JC\MD	Ok
2	VSTDCCC050	VSTDCCC050	VN088437.D	04 Dec 2025 11:44	pH#Lot#V12668	JC\MD	Ok,M
3	VN1204MBL01	VN1204MBL01	VN088438.D	04 Dec 2025 12:17		JC\MD	Ok
4	VN1204WBL01	VN1204WBL01	VN088439.D	04 Dec 2025 12:38		JC\MD	Ok
5	VN1204MBS01	VN1204MBS01	VN088440.D	04 Dec 2025 13:11		JC\MD	Ok,M
6	Q3768-01	20071	VN088441.D	04 Dec 2025 13:31	Need Soil Run	JC\MD	Not Ok
7	IBLK	IBLK	VN088442.D	04 Dec 2025 13:52		JC\MD	Ok
8	VN1204WBS01	VN1204WBS01	VN088443.D	04 Dec 2025 14:12		JC\MD	Ok,M
9	VN1204WBSD01	VN1204WBSD01	VN088444.D	04 Dec 2025 14:45		JC\MD	Ok,M
10	Q3756-03	BR-01-190-120225	VN088445.D	04 Dec 2025 15:06	vial A pH<2	JC\MD	Ok
11	Q3756-01	OW-04B-26.5-120225	VN088446.D	04 Dec 2025 15:26	vial A pH<2	JC\MD	Ok
12	Q3756-05	MW-04-12-120225	VN088447.D	04 Dec 2025 15:47	vial A pH<2	JC\MD	Ok
13	Q3756-07	MW-13B-47.5-120225	VN088448.D	04 Dec 2025 16:07	vial A pH<2	JC\MD	Ok
14	Q3757-12	TB01-120325	VN088449.D	04 Dec 2025 16:28	vial A pH<2 TB	JC\MD	Ok
15	Q3757-11	EB01-120325	VN088450.D	04 Dec 2025 16:48	vial A pH<2 EB	JC\MD	Ok,M
16	Q3757-01	MW-14A-17.5-120325	VN088451.D	04 Dec 2025 17:09	Need Straight Run	JC\MD	Not Ok
17	Q3757-03	BR-03D-490-120325	VN088452.D	04 Dec 2025 17:29	Need Straight Run	JC\MD	Not Ok
18	Q3757-05	MW-14B-46-120325	VN088453.D	04 Dec 2025 17:50	Need Straight Run	JC\MD	Not Ok

Instrument ID: MSVOA_N

Daily Analysis Runlog For Sequence/QC Batch ID # VN120425

Review By	John Carlone	Review On	12/5/2025 8:23:32 AM
Supervise By	Semsettin Yesilyurt	Supervise On	12/5/2025 2:14:29 PM
SubDirectory	VN120425	HP Acquire Method	HP Processing Method 82N112425W.M
STD. NAME	STD REF.#		
Tune/Reschk Initial Calibration Stds	VP136505		
CCC Internal Standard/PEM ICV/I.BLK Surrogate Standard MS/MSD Standard LCS Standard	VP136506,VP136507		

19	Q3757-07	RMW-05B-90.5-120325	VN088454.D	04 Dec 2025 18:10	Need Straight Run	JC\MD	Not Ok
20	Q3757-09	OW-05B-72-120325	VN088455.D	04 Dec 2025 18:31	Need Straight Run	JC\MD	Not Ok
21	Q3750-07	FENCE WC-1	VN088456.D	04 Dec 2025 18:51	vial B pH#5.0	JC\MD	Ok
22	Q3750-08	FENCE WC-2	VN088457.D	04 Dec 2025 19:11	vial B pH#5.0	JC\MD	Ok
23	PB170809TB	PB170809TB	VN088458.D	04 Dec 2025 19:32		JC\MD	Ok
24	Q3775-01	NGKO	VN088459.D	04 Dec 2025 19:52	Need Straight Run	JC\MD	Not Ok
25	IBLK	IBLK	VN088460.D	04 Dec 2025 20:13		JC\MD	Ok
26	VSTDCCC050	VSTDCCC050EC	VN088461.D	04 Dec 2025 20:33		JC\MD	Ok,M

M : Manual Integration

Instrument ID: MSVOA_X

Daily Analysis Runlog For Sequence/QC Batch ID # VX120425

Review By	John Carlone	Review On	12/5/2025 8:05:59 AM
Supervise By	Semsettin Yesilyurt	Supervise On	12/5/2025 2:14:46 PM
SubDirectory	VX120425	HP Acquire Method	HP Processing Method 82X120425

STD. NAME	STD REF.#
Tune/Reschk	VP136497
Initial Calibration Stds	VP136526,VP136527,VP136528,VP136529,VP136530,VP136531
CCC	
Internal Standard/PEM	
ICV/I.BLK	VP136532
Surrogate Standard	
MS/MSD Standard	
LCS Standard	

Sr#	Sampled	ClientID	Data File Name	Date-Time	Comment	Operator	Status
1	BFB	BFB	VX048713.D	04 Dec 2025 08:15		JC/MD	Ok
2	VSTDICC001	VSTDICC001	VX048714.D	04 Dec 2025 08:47		JC/MD	Ok,M
3	VSTDICC005	VSTDICC005	VX048715.D	04 Dec 2025 09:16		JC/MD	Ok,M
4	VSTDICC020	VSTDICC020	VX048716.D	04 Dec 2025 09:36	Comp. #13 on Linear Regression	JC/MD	Ok,M
5	VSTDICCC050	VSTDICCC050	VX048717.D	04 Dec 2025 09:57	Comp. #95 on Quadratic Regression	JC/MD	Ok,M
6	VSTDICC100	VSTDICC100	VX048718.D	04 Dec 2025 10:17		JC/MD	Ok,M
7	VSTDICC150	VSTDICC150	VX048719.D	04 Dec 2025 10:38		JC/MD	Ok,M
8	IBLK	IBLK	VX048720.D	04 Dec 2025 10:58		JC/MD	Ok
9	VSTDICV050	ICVVX120425	VX048721.D	04 Dec 2025 14:04		JC/MD	Ok,M

M : Manual Integration

Instrument ID: MSVOA_X

Daily Analysis Runlog For Sequence/QC Batch ID # VX120525

Review By	John Carlone	Review On	12/8/2025 8:24:07 AM
Supervise By	Semsettin Yesilyurt	Supervise On	12/8/2025 11:02:48 AM
SubDirectory	VX120525	HP Acquire Method	HP Processing Method 82X120425

STD. NAME	STD REF.#
Tune/Reschk Initial Calibration Stds	VP136523
CCC Internal Standard/PEM ICV/I.BLK Surrogate Standard MS/MSD Standard LCS Standard	VP136524,VP136525

Sr#	SampleID	ClientID	Data File Name	Date-Time	Comment	Operator	Status
1	BFB	BFB	VX048722.D	05 Dec 2025 08:14		JC/MD	Ok
2	VSTDCCC050	VSTDCCC050	VX048723.D	05 Dec 2025 08:45	pH#Lot#V12668	JC/MD	Ok,M
3	VX1205MBL01	VX1205MBL01	VX048724.D	05 Dec 2025 09:14		JC/MD	Ok
4	VX1205WBL01	VX1205WBL01	VX048725.D	05 Dec 2025 10:34		JC/MD	Ok
5	VX1205WBS01	VX1205WBS01	VX048726.D	05 Dec 2025 11:31		JC/MD	Ok,M
6	VX1205WBSD01	VX1205WBSD01	VX048727.D	05 Dec 2025 11:59		JC/MD	Ok,M
7	Q3775-01	NGKO	VX048728.D	05 Dec 2025 12:20	vial A pH#5.0 Surrogate fail	JC/MD	ReRun
8	Q3775-01RE	NGKORE	VX048729.D	05 Dec 2025 12:50	vial B pH#5.0 Surrogate fail	JC/MD	Confirms
9	IBLK	IBLK	VX048730.D	05 Dec 2025 13:10		JC/MD	Ok
10	Q3757-01	MW-14A-17.5-120325	VX048731.D	05 Dec 2025 13:32	vial A pH<2	JC/MD	Ok
11	Q3757-03	BR-03D-490-120325	VX048732.D	05 Dec 2025 13:53	vial A pH<2	JC/MD	Ok
12	Q3757-05	MW-14B-46-120325	VX048733.D	05 Dec 2025 14:13	vial A pH<2 Surrogate fail	JC/MD	ReRun
13	Q3757-07	RMW-05B-90.5-120325	VX048734.D	05 Dec 2025 14:33	vial A pH<2	JC/MD	Ok
14	Q3757-09	OW-05B-72-120325	VX048735.D	05 Dec 2025 14:54	vial A pH<2 Surrogate fail	JC/MD	ReRun
15	Q3789-01	STORAGE-BLANK-SO	VX048736.D	05 Dec 2025 15:14	vial A pH<2	JC/MD	Ok
16	Q3789-02	STORAGE-BLANK-WA	VX048737.D	05 Dec 2025 15:35	vial A pH<2	JC/MD	Ok
17	Q3789-03	STORAGE-BLANK-WA	VX048738.D	05 Dec 2025 15:55	vial A pH<2 Surrogate fail	JC/MD	ReRun
18	Q3789-04	STORAGE-BLANK-SAI	VX048739.D	05 Dec 2025 16:15	vial A pH<2 Surrogate fail	JC/MD	ReRun

Instrument ID: MSVOA_X

Daily Analysis Runlog For Sequence/QC Batch ID # VX120525

Review By	John Carlone	Review On	12/8/2025 8:24:07 AM
Supervise By	Semsettin Yesilyurt	Supervise On	12/8/2025 11:02:48 AM
SubDirectory	VX120525	HP Acquire Method	HP Processing Method 82X120425

STD. NAME	STD REF.#
Tune/Reschk Initial Calibration Stds	VP136523
CCC Internal Standard/PEM ICV/I.BLK Surrogate Standard MS/MSD Standard LCS Standard	VP136524,VP136525

19	VSTDCCC050	VSTDCCC050EC	VX048740.D	05 Dec 2025 16:36		JC/MD	Ok,M
----	------------	--------------	------------	-------------------	--	-------	------

M : Manual Integration

Instrument ID: MSVOA_X

Daily Analysis Runlog For Sequence/QC Batch ID # VX120825

Review By	John Carlone	Review On	12/9/2025 8:11:51 AM
Supervise By	Maresh Dadoda	Supervise On	12/9/2025 12:45:16 PM
SubDirectory	VX120825	HP Acquire Method	HP Processing Method 82X120425W.M

STD. NAME	STD REF.#
Tune/Reschk Initial Calibration Stds	VP136561
CCC Internal Standard/PEM ICV/I.BLK Surrogate Standard MS/MSD Standard LCS Standard	VP136570,VP136571

Sr#	SampleID	ClientID	Data File Name	Date-Time	Comment	Operator	Status
1	BFB	BFB	VX048741.D	08 Dec 2025 08:17		JC/MD	Ok
2	VSTDCCC050	VSTDCCC050	VX048742.D	08 Dec 2025 09:20	pH#Lot#V12668	JC/MD	Ok,M
3	VX1208MBL01	VX1208MBL01	VX048743.D	08 Dec 2025 09:49		JC/MD	Ok
4	VX1208WBL01	VX1208WBL01	VX048744.D	08 Dec 2025 11:15		JC/MD	Ok
5	VX1208WBS01	VX1208WBS01	VX048745.D	08 Dec 2025 11:44		JC/MD	Ok,M
6	VX1208WBSD01	VX1208WBSD01	VX048746.D	08 Dec 2025 12:12		JC/MD	Ok,M
7	Q3757-05RE	MW-14B-46-120325RE	VX048747.D	08 Dec 2025 12:33	vial B pH<2 Surrogate fail	JC/MD	Confirms
8	Q3757-09RE	OW-05B-72-120325RE	VX048748.D	08 Dec 2025 12:53	vial B pH<2 Surrogate fail	JC/MD	Confirms
9	Q3787-01	MW-15B-42.5-120425	VX048749.D	08 Dec 2025 13:14	vial A pH<2	JC/MD	Ok
10	Q3787-02MS	MW-15B-42.5-120425	VX048750.D	08 Dec 2025 13:35	Not Spiked	JC/MD	Not Ok
11	Q3787-03MSD	MW-15B-42.5-120425	VX048751.D	08 Dec 2025 13:55	Not Spiked	JC/MD	Not Ok
12	Q3787-07	OWBR-02-170-120425	VX048752.D	08 Dec 2025 14:16	vial A pH<2 Surrogate fail	JC/MD	ReRun
13	Q3787-09	OWBR-02-170-120425	VX048753.D	08 Dec 2025 14:36	vial A pH<2 Surrogate fail	JC/MD	ReRun
14	Q3787-11	OW-03B-51.5-120425	VX048754.D	08 Dec 2025 14:57	vial A pH<2 Surrogate fail	JC/MD	ReRun
15	Q3787-13	OW-03B-51.5-120425	VX048755.D	08 Dec 2025 15:18	vial A pH<2 Surrogate fail	JC/MD	ReRun
16	Q3787-15	OW-08B-72.5-120425	VX048756.D	08 Dec 2025 15:38	vial A pH<2	JC/MD	Ok
17	Q3787-17	OW-08B-72.5-120425	VX048757.D	08 Dec 2025 15:59	vial A pH<2 Surrogate fail	JC/MD	ReRun
18	Q3787-19	OW-02B-21.2-120425	VX048758.D	08 Dec 2025 16:19	vial A pH<2 Surrogate fail	JC/MD	ReRun

Instrument ID: MSVOA_X

Daily Analysis Runlog For Sequence/QC Batch ID # VX120825

Review By	John Carlone	Review On	12/9/2025 8:11:51 AM
Supervise By	Maresh Dadoda	Supervise On	12/9/2025 12:45:16 PM
SubDirectory	VX120825	HP Acquire Method	HP Processing Method 82X120425W.M
STD. NAME	STD REF.#		
Tune/Reschk Initial Calibration Stds	VP136561		
CCC Internal Standard/PEM ICV/I.BLK Surrogate Standard MS/MSD Standard LCS Standard	VP136570,VP136571		

19	Q3787-21	TB01-120425	VX048759.D	08 Dec 2025 16:40	vial A pH<2 TB;Surrogate fail	JC/MD	ReRun
20	Q3788-07	OW-01B-66.5-120525	VX048760.D	08 Dec 2025 17:00	vial A pH<2 Surrogate fail	JC/MD	ReRun
21	Q3788-09	TA-BR-05-465-120525	VX048761.D	08 Dec 2025 17:21	vial A pH<2 Surrogate fail	JC/MD	ReRun
22	Q3788-11	MW-16B-87.5-120525	VX048762.D	08 Dec 2025 17:41	vial A pH<2 Surrogate fail	JC/MD	ReRun
23	Q3788-13	TB01-120525	VX048763.D	08 Dec 2025 18:02	vial A pH<2 TB	JC/MD	Ok
24	Q3787-02MS	MW-15B-42.5-120425	VX048764.D	08 Dec 2025 18:22	vial A pH<2	JC/MD	Ok,M
25	Q3787-03MSD	MW-15B-42.5-120425	VX048765.D	08 Dec 2025 18:43	vial A pH<2	JC/MD	Ok,M
26	VSTDCCC050	VSTDCCC050EC	VX048766.D	08 Dec 2025 19:03		JC/MD	Ok,M

M : Manual Integration



SHIPPING DOCUMENTS

ALLIANCE PROJECT NO.	
QUOTE NO.	Q3757
COC Number	2047320

CLIENT INFORMATION

REPORT TO BE SENT TO:

COMPANY: Jacobs
ADDRESS: 412 Mt Kumbie Ave Suite H100
CITY Morristown STATE: NJ ZIP: 07960
ATTENTION: John Yubante
PHONE: (281) 414-1719 FAX:

CLIENT PROJECT INFORMATION

PROJECT NAME: STC PTC
PROJECT NO.: D4033126 LOCATION: Princeton Junction
PROJECT MANAGER: Mary Murphy
e-mail: Mary.Murphy@Jacobs.com
PHONE: (201) 936-0586 FAX:

CLIENT BILLING INFORMATION

BILL TO: Mary Murphy PO#:
ADDRESS:
CITY STATE: ZIP:
ATTENTION: PHONE:

ANALYSIS

DATA TURNAROUND INFORMATION

FAX (RUSH) Standard DAYS*
HARDCOPY (DATA PACKAGE): DAYS*
EDD: DAYS*
*TO BE APPROVED BY CHEMTECH
STANDARD HARDCOPY TURNAROUND TIME IS 10 BUSINESS

DATA DELIVERABLE INFORMATION

☐ Level 1 (Results Only) ☐ Level 4 (QC + Full Raw Data)
☐ Level 2 (Results + QC) ☐ NJ Reduced ☐ US EPA CLP
☒ Level 3 (Results + QC) ☐ NYS ASP A ☐ NYS ASP B
+ Raw Data ☐ Other
☐ EDD FORMAT

*Site Specific VOCs
(82100-51M)
Time VOCs Site Specific
(SFAM) 114-114-114
114-114-114
(82100-51M)*

PRESERVATIVES

COMMENTS

ALLIANCE SAMPLE ID	PROJECT SAMPLE IDENTIFICATION	SAMPLE MATRIX	SAMPLE TYPE		SAMPLE COLLECTION		# OF BOTTLES	PRESERVATIVES									COMMENTS ← Specify Preservatives A-HCl D-NaOH B-HNO3 E-ICE C-H2SO4 F-OTHER
			COMP	GRAB	DATE	TIME		A/E	A/E	E							
1.	EB01-120325	DI		X	12/3/25	1530	8	X	X	X							
2.	TB01-120325	DI		X	12/3/25	0800	4	X	X								
3.																	
4.																	
5.																	
6.																	
7.																	
8.																	
9.																	
10.																	

SAMPLE CUSTODY MUST BE DOCUMENTED BELOW EACH TIME SAMPLES CHANGE POSSESSION INCLUDING COURIER DELIVERY

RELINQUISHED BY SAMPLER: 1. <u>[Signature]</u>	DATE/TIME: <u>12-3-25 1610</u>	RECEIVED BY: <u>[Signature]</u> <u>1610</u>	Conditions of bottles or coolers at receipt: ☐ COMPLIANT ☐ NON COMPLIANT ☐ COOLER TEMP <u>2.8</u> °C
RELINQUISHED BY SAMPLER: 2. <u>[Signature]</u>	DATE/TIME:	RECEIVED BY: <u>[Signature]</u>	Comments: <u>See work order for list of site specific VOCs</u>
RELINQUISHED BY SAMPLER: 3. <u>[Signature]</u>	DATE/TIME: <u>12-3-25 1747</u>	RECEIVED BY: <u>[Signature]</u>	

Page 2 of 2 CLIENT: ☐ Hand Delivered ☐ Other Shipment Complete ☐ YES ☐ NO

Laboratory Certification

Certified By	License No.
Connecticut	PH-0830
DOD ELAP (ANAB)	L2219
Maine	2024021
Maryland	296
New Hampshire	255425
New Jersey	20012
New York	11376
Pennsylvania	68-00548
Soil Permit	525-24-234-08441
Texas	TX-C25-00189
Virginia	460312

LOGIN REPORT/SAMPLE TRANSFER

Order ID : Q3757	JACO05	Order Date : 12/3/2025 9:25:00 AM	Project Mgr : Deepak
Client Name : JACOBS Engineering Grou		Project Name : Former Schlumberger STC	Report Type : Level 3
Client Contact : John Ynfante		Receive DateTime : 12/3/2025 5:47:00 PM	EDD Type : CH2MHILL
Invoice Name : JACOBS Engineering Grou		Purchase Order :	Hard Copy Date :
Invoice Contact : John Ynfante			Date Signoff : 12/4/2025 2:28:49 PM

LAB ID	CLIENT ID	MATRIX	SAMPLE DATE	SAMPLE TIME	TEST	TEST GROUP	METHOD	FAX DATE	DUE DATES
Q3757-01	MW-14A-17.5-120325	Water	12/03/2025	10:20					
					VOCMS Group3		8260-Low	10 Bus. Days	
Q3757-02	MW-14A-17.5-120325-SIM	Water	12/03/2025	10:20					
					VOC-SIM		SFAM_VOCSIM	10 Bus. Days	
Q3757-03	BR-03D-490-120325	Water	12/03/2025	12:50					
					VOCMS Group3		8260-Low	10 Bus. Days	
Q3757-04	BR-03D-490-120325-SIM	Water	12/03/2025	12:50					
					VOC-SIM		SFAM_VOCSIM	10 Bus. Days	
Q3757-05	MW-14B-46-120325	Water	12/03/2025	13:00					
					VOCMS Group3		8260-Low	10 Bus. Days	
Q3757-06	MW-14B-46-120325-SIM	Water	12/03/2025	13:00					
					VOC-SIM		SFAM_VOCSIM	10 Bus. Days	
Q3757-07	RMW-05B-90.5-120325	Water	12/03/2025	13:50					
					VOCMS Group3		8260-Low	10 Bus. Days	
Q3757-08	RMW-05B-90.5-120325-SIM	Water	12/03/2025	13:50					



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

LOGIN REPORT/SAMPLE TRANSFER

Order ID : Q3757	JACO05	Order Date : 12/3/2025 9:25:00 AM	Project Mgr : Deepak
Client Name : JACOBS Engineering Grou		Project Name : Former Schlumberger STC	Report Type : Level 3
Client Contact : John Ynfante		Receive DateTime : 12/3/2025 5:47:00 PM	EDD Type : CH2MHILL
Invoice Name : JACOBS Engineering Grou		Purchase Order :	Hard Copy Date :
Invoice Contact : John Ynfante			Date Signoff : 12/4/2025 2:28:49 PM

LAB ID	CLIENT ID	MATRIX	SAMPLE DATE	SAMPLE TIME	TEST	TEST GROUP	METHOD	FAX DATE	DUE DATES
					VOC-SIM		SFAM_VOCSIM	10 Bus. Days	
Q3757-09	OW-05B-72-120325	Water	12/03/2025	14:40					
					VOCMS Group3		8260-Low	10 Bus. Days	
Q3757-10	OW-05B-72-120325-SIM	Water	12/03/2025	14:40					
					VOC-SIM		SFAM_VOCSIM	10 Bus. Days	
Q3757-11	EB01-120325	Water	12/03/2025	15:30					
					VOC-SIM		SFAM_VOCSIM	10 Bus. Days	
					VOCMS Group3		8260-Low	10 Bus. Days	
Q3757-12	TB01-120325	Water	12/03/2025	08:00					
					VOC-SIM		SFAM_VOCSIM	10 Bus. Days	
					VOCMS Group3		8260-Low	10 Bus. Days	
Q3757-14	VHBLK001	Water	12/03/2025	17:47					
					VOC-SIM		SFAM_VOCSIM	10 Bus. Days	



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

LOGIN REPORT/SAMPLE TRANSFER

Order ID : Q3757	JACO05	Order Date : 12/3/2025 9:25:00 AM	Project Mgr : Deepak
Client Name : JACOBS Engineering Grou		Project Name : Former Schlumberger STC	Report Type : Level 3
Client Contact : John Ynfante		Receive DateTime : 12/3/2025 5:47:00 PM	EDD Type : CH2MHILL
Invoice Name : JACOBS Engineering Grou		Purchase Order :	Hard Copy Date :
Invoice Contact : John Ynfante			Date Signoff : 12/4/2025 2:28:49 PM

LAB ID	CLIENT ID	MATRIX	SAMPLE DATE	SAMPLE TIME	TEST	TEST GROUP	METHOD	FAX DATE	DUE DATES
--------	-----------	--------	----------------	----------------	------	------------	--------	----------	--------------

Relinquished By : ap
Date / Time : 12/4/25 9:30

Received By : JC
Date / Time : 12/4/25 9:30
Storage Area : VOA Refridgerator Room