

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:	Q3743	OrderDate:	12/1/2025 9:45:00 AM
Client:	Remington & Vernick	Project:	Saddler Property
Contact:	Justin Zarzecki	Location:	A12,VOA Lab

LabID	ClientID	Matrix	Test	Method	Sample Date	Prep Date	Anal Date	Received
Q3743-01	TW-1125-1	Water	VOC-SIM	SFAM_VOCSI M	11/26/25		12/03/25	11/26/25
			VOC-TCLVOA-10	8260-Low			12/01/25	
Q3743-02	TW-1125-2	Water	VOC-SIM	SFAM_VOCSI M	11/26/25		12/03/25	11/26/25
			VOC-TCLVOA-10	8260-Low			12/01/25	
Q3743-03	TB	Water	VOC-SIM	SFAM_VOCSI M	11/26/25		12/03/25	11/26/25
			VOC-TCLVOA-10	8260-Low			12/01/25	
Q3743-04	TB	Water	VOC-SIM	SFAM_VOCSI M	11/26/25		12/03/25	11/26/25
			VOC-TCLVOA-10	8260-Low			12/01/25	
Q3743-05	TB	Water	VOC-SIM	SFAM_VOCSI M	11/26/25		12/03/25	11/26/25
			VOC-TCLVOA-10	8260-Low			12/01/25	

Hit Summary Sheet
SW-846

SDG No.: Q3743

Client: Remington & Vernick

Sample ID	Client ID	Matrix	Parameter	Concentration	C	MDL	RDL	Units
Client ID:	TW-1125-1							
Q3743-01	TW-1125-1	Water	Acetone	21.6		1.50	5.00	ug/L
			Total Voc :	21.6				
			Total Concentration:	21.6				
Client ID:	TW-1125-2							
Q3743-02	TW-1125-2	Water	Acetone	19.0		1.50	5.00	ug/L
			Total Voc :	19.0				
			Total Concentration:	19.0				



QC SUMMARY

Surrogate Summary

SDG No.: **Q3743**

Client: **Remington & Vernick**

Analytical Method: **SW8260-Low**

Lab Sample ID	Client ID	Parameter	Spike	Result	Recovery (%)	Qual	Limits (%)	
							Low	High
Q3743-01	TW-1125-1	1,2-Dichloroethane-d4	50	61.7	123		74	125
		Dibromofluoromethane	50	53.8	108		75	124
		Toluene-d8	50	50.2	100		86	113
		4-Bromofluorobenzene	50	53.4	107		77	121
Q3743-02	TW-1125-2	1,2-Dichloroethane-d4	50	61.2	122		74	125
		Dibromofluoromethane	50	53.0	106		75	124
		Toluene-d8	50	50.1	100		86	113
		4-Bromofluorobenzene	50	53.5	107		77	121
Q3743-03	TB	1,2-Dichloroethane-d4	50	60.1	120		74	125
		Dibromofluoromethane	50	52.4	105		75	124
		Toluene-d8	50	50.1	100		86	113
		4-Bromofluorobenzene	50	53.5	107		77	121
Q3743-04		1,2-Dichloroethane-d4	50	60.9	122		74	125
		Dibromofluoromethane	50	53.0	106		75	124
		Toluene-d8	50	49.8	100		86	113
		4-Bromofluorobenzene	50	53.2	106		77	121
VN1201WBL01	VN1201WBL01	1,2-Dichloroethane-d4	50	57.8	116		74	125
		Dibromofluoromethane	50	52.2	104		75	124
		Toluene-d8	50	50.0	100		86	113
		4-Bromofluorobenzene	50	52.4	105		77	121
VN1201WBS01	VN1201WBS01	1,2-Dichloroethane-d4	50	56.0	112		74	125
		Dibromofluoromethane	50	55.1	110		75	124
		Toluene-d8	50	54.9	110		86	113
		4-Bromofluorobenzene	50	53.3	107		77	121
VN1201WBSD0	VN1201WBSD01	1,2-Dichloroethane-d4	50	52.2	104		74	125
		Dibromofluoromethane	50	52.3	105		75	124
		Toluene-d8	50	53.2	106		86	113
		4-Bromofluorobenzene	50	51.1	102		77	121

Laboratory Control Sample/Laboratory Control Sample Duplicate Summary

SW-846

SDG No.:	<u>Q3743</u>	Analytical Method:	<u>SW8260-Low</u>
Client:	<u>Remington & Vernick</u>	Datafile :	<u>VN088373.D</u>

Lab Sample ID	Parameter	Spike	Result	Unit	Rec	RPD	Qual	Low	Limits High	RPD
VN1201WBS01	Dichlorodifluoromethane	20	18.2	ug/L	91			69	116	
	Chloromethane	20	19.4	ug/L	97			65	116	
	Vinyl chloride	20	18.9	ug/L	95			65	117	
	Bromomethane	20	19.7	ug/L	99			58	125	
	Chloroethane	20	19.5	ug/L	98			56	128	
	Trichlorofluoromethane	20	18.6	ug/L	93			73	115	
	1,1,2-Trichlorotrifluoroethane	20	18.6	ug/L	93			80	112	
	1,1-Dichloroethene	20	19.0	ug/L	95			74	110	
	Acetone	100	87.9	ug/L	88			60	125	
	Carbon disulfide	20	19.4	ug/L	97			64	112	
	Methyl tert-butyl Ether	20	19.9	ug/L	100			78	114	
	Methyl Acetate	20	19.1	ug/L	96			57	150	
	Methylene Chloride	20	20.0	ug/L	100			72	114	
	trans-1,2-Dichloroethene	20	19.6	ug/L	98			75	108	
	1,1-Dichloroethane	20	19.8	ug/L	99			78	112	
	Cyclohexane	20	17.9	ug/L	90			75	110	
	2-Butanone	100	96.0	ug/L	96			65	122	
	Carbon Tetrachloride	20	18.3	ug/L	92			77	113	
	cis-1,2-Dichloroethene	20	20.1	ug/L	101			77	110	
	Bromochloromethane	20	19.3	ug/L	97			70	124	
	Chloroform	20	20.3	ug/L	102			79	113	
	1,1,1-Trichloroethane	20	18.8	ug/L	94			80	108	
	Methylcyclohexane	20	17.1	ug/L	86			72	115	
	Benzene	20	19.4	ug/L	97			82	109	
	1,2-Dichloroethane	20	19.6	ug/L	98			80	115	
	Trichloroethene	20	18.8	ug/L	94			77	113	
	1,2-Dichloropropane	20	19.8	ug/L	99			83	111	
	Bromodichloromethane	20	19.5	ug/L	98			83	110	
	4-Methyl-2-Pentanone	100	98.3	ug/L	98			74	118	
	Toluene	20	19.5	ug/L	98			82	110	
	t-1,3-Dichloropropene	20	19.9	ug/L	100			79	110	
	cis-1,3-Dichloropropene	20	20.1	ug/L	101			82	110	
	1,1,2-Trichloroethane	20	20.2	ug/L	101			83	112	
	2-Hexanone	100	93.5	ug/L	94			73	117	
	Dibromochloromethane	20	20.1	ug/L	101			82	110	
	1,2-Dibromoethane	20	20.6	ug/L	103			81	110	
	Tetrachloroethene	20	17.9	ug/L	90			67	123	
	Chlorobenzene	20	19.5	ug/L	98			82	109	
	Ethyl Benzene	20	18.6	ug/L	93			83	109	
	m/p-Xylenes	40	37.0	ug/L	93			82	110	
	o-Xylene	20	18.3	ug/L	92			83	109	
	Styrene	20	19.6	ug/L	98			80	111	
	Bromoform	20	20.0	ug/L	100			79	109	
	Isopropylbenzene	20	18.6	ug/L	93			83	112	
	1,1,2,2-Tetrachloroethane	20	20.1	ug/L	101			76	118	
	1,3-Dichlorobenzene	20	19.0	ug/L	95			82	108	
	1,4-Dichlorobenzene	20	18.5	ug/L	93			82	107	
	1,2-Dichlorobenzene	20	19.4	ug/L	97			82	109	

Laboratory Control Sample/Laboratory Control Sample Duplicate Summary

SW-846

SDG No.: Q3743

Analytical Method: SW8260-Low

Client: Remington & Vernick

Datafile : VN088373.D

Lab Sample ID	Parameter	Spike	Result	Unit	Rec	RPD	Qual	Low	Limits	RPD
									High	
VN1201WBS01	1,2-Dibromo-3-Chloropropane	20	18.2	ug/L	91			68	112	
	1,2,4-Trichlorobenzene	20	18.5	ug/L	93			75	113	
	1,2,3-Trichlorobenzene	20	18.3	ug/L	92			76	114	

Laboratory Control Sample/Laboratory Control Sample Duplicate Summary

SW-846

SDG No.:	<u>Q3743</u>	Analytical Method:	<u>SW8260-Low</u>
Client:	<u>Remington & Vernick</u>	Datafile :	<u>VN088374.D</u>

Lab Sample ID	Parameter	Spike	Result	Unit	Rec	RPD	Qual	Low	Limits High	RPD
VN1201WBSD01	Dichlorodifluoromethane	20	18.9	ug/L	95	4		69	116	19
	Chloromethane	20	19.4	ug/L	97	0		65	116	21
	Vinyl chloride	20	18.7	ug/L	94	1		65	117	19
	Bromomethane	20	20.2	ug/L	101	2		58	125	30
	Chloroethane	20	19.3	ug/L	97	1		56	128	20
	Trichlorofluoromethane	20	18.7	ug/L	94	1		73	115	16
	1,1,2-Trichlorotrifluoroethane	20	18.5	ug/L	93	0		80	112	20
	1,1-Dichloroethene	20	18.8	ug/L	94	1		74	110	20
	Acetone	100	87.0	ug/L	87	1		60	125	27
	Carbon disulfide	20	18.8	ug/L	94	3		64	112	17
	Methyl tert-butyl Ether	20	19.4	ug/L	97	3		78	114	20
	Methyl Acetate	20	19.0	ug/L	95	1		57	150	20
	Methylene Chloride	20	20.2	ug/L	101	1		72	114	20
	trans-1,2-Dichloroethene	20	19.6	ug/L	98	0		75	108	16
	1,1-Dichloroethane	20	19.1	ug/L	96	3		78	112	20
	Cyclohexane	20	18.5	ug/L	93	3		75	110	20
	2-Butanone	100	96.2	ug/L	96	0		65	122	26
	Carbon Tetrachloride	20	18.8	ug/L	94	2		77	113	15
	cis-1,2-Dichloroethene	20	19.4	ug/L	97	4		77	110	20
	Bromochloromethane	20	18.5	ug/L	93	4		70	124	20
	Chloroform	20	19.6	ug/L	98	4		79	113	20
	1,1,1-Trichloroethane	20	19.0	ug/L	95	1		80	108	20
	Methylcyclohexane	20	18.5	ug/L	93	8		72	115	20
	Benzene	20	19.1	ug/L	96	1		82	109	15
	1,2-Dichloroethane	20	19.4	ug/L	97	1		80	115	20
	Trichloroethene	20	19.1	ug/L	96	2		77	113	15
	1,2-Dichloropropane	20	19.2	ug/L	96	3		83	111	21
	Bromodichloromethane	20	19.3	ug/L	97	1		83	110	21
	4-Methyl-2-Pentanone	100	98.5	ug/L	99	1		74	118	25
	Toluene	20	19.8	ug/L	99	1		82	110	16
	t-1,3-Dichloropropene	20	19.1	ug/L	96	4		79	110	20
	cis-1,3-Dichloropropene	20	19.5	ug/L	98	3		82	110	16
	1,1,2-Trichloroethane	20	20.0	ug/L	100	1		83	112	20
	2-Hexanone	100	94.4	ug/L	94	0		73	117	25
	Dibromochloromethane	20	20.3	ug/L	102	1		82	110	20
	1,2-Dibromoethane	20	19.9	ug/L	100	3		81	110	20
	Tetrachloroethene	20	19.0	ug/L	95	5		67	123	15
	Chlorobenzene	20	19.8	ug/L	99	1		82	109	15
	Ethyl Benzene	20	18.9	ug/L	95	2		83	109	16
	m/p-Xylenes	40	37.8	ug/L	95	2		82	110	20
	o-Xylene	20	19.2	ug/L	96	4		83	109	20
	Styrene	20	19.4	ug/L	97	1		80	111	17
	Bromoform	20	20.1	ug/L	101	1		79	109	20
	Isopropylbenzene	20	19.3	ug/L	97	4		83	112	22
	1,1,2,2-Tetrachloroethane	20	19.5	ug/L	98	3		76	118	20
	1,3-Dichlorobenzene	20	19.5	ug/L	98	3		82	108	20
	1,4-Dichlorobenzene	20	19.0	ug/L	95	2		82	107	20
	1,2-Dichlorobenzene	20	19.0	ug/L	95	2		82	109	20

Laboratory Control Sample/Laboratory Control Sample Duplicate Summary

SW-846

SDG No.: Q3743

Analytical Method: SW8260-Low

Client: Remington & Vernick

Datafile : VN088374.D

Lab Sample ID	Parameter	Spike	Result	Unit	Rec	RPD	Qual	Low	Limits	RPD
									High	
VN1201WBSD01	1,2-Dibromo-3-Chloropropane	20	18.5	ug/L	93	2		68	112	26
	1,2,4-Trichlorobenzene	20	18.0	ug/L	90	3		75	113	21
	1,2,3-Trichlorobenzene	20	19.1	ug/L	96	4		76	114	22

VOLATILE METHOD BLANK SUMMARY

Client ID

VN1201WBL01

Lab Name: Alliance

Contract: REMI01

Lab Code: ACE

SDG NO.: Q3743

Lab File ID: VN088372.D

Lab Sample ID: VN1201WBL01

Date Analyzed: 12/01/2025

Time Analyzed: 10:53

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
VN1201WBS01	VN1201WBS01	VN088373.D	12/01/2025
VN1201WBSD01	VN1201WBSD01	VN088374.D	12/01/2025
TB	Q3743-03	VN088389.D	12/01/2025
TB	Q3743-04	VN088390.D	12/01/2025
TW-1125-1	Q3743-01	VN088392.D	12/01/2025
TW-1125-2	Q3743-02	VN088393.D	12/01/2025

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: REMI01

Lab Code: ACE

SDG NO.: Q3743

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,
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VOLATILE ORGANIC INSTRUMENT PERFORMANCE CHECK
BROMOFLUOROBENZENE (BFB)

Lab Name: Alliance

Contract: REMI01

Lab Code: ACE

SDG NO.: Q3743

Lab File ID: VN088369.D

BFB Injection Date: 12/01/2025

Instrument ID: MSVOA_N

BFB Injection Time: 08:49

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	22.9
75	30.0 - 60.0% of mass 95	59.1
95	Base Peak, 100% relative abundance	100
96	5.0 - 9.0% of mass 95	6.7
173	Less than 2.0% of mass 174	0.9 (1.2) 1
174	50.0 - 100.0% of mass 95	71
175	5.0 - 9.0% of mass 174	5.7 (8) 1
176	95.0 - 101.0% of mass 174	68.9 (97.1) 1
177	5.0 - 9.0% of mass 176	5.2 (7.5) 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	VN088370.D	12/01/2025	09:58
VN1201WBL01	VN1201WBL01	VN088372.D	12/01/2025	10:53
VN1201WBS01	VN1201WBS01	VN088373.D	12/01/2025	11:27
VN1201WBSD01	VN1201WBSD01	VN088374.D	12/01/2025	12:10
TB	Q3743-03	VN088389.D	12/01/2025	17:18
TB	Q3743-04	VN088390.D	12/01/2025	17:39
TW-1125-1	Q3743-01	VN088392.D	12/01/2025	18:20
TW-1125-2	Q3743-02	VN088393.D	12/01/2025	18:41

VOLATILE INTERNAL STANDARD AREA AND RT SUMMARY

Lab Name: Alliance Contract: REMI01
 Lab Code: ACE SDG NO.: Q3743
 Lab File ID: VN088370.D Date Analyzed: 12/01/2025
 Instrument ID: MSVOA_N Time Analyzed: 09:58
 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	262500	8.20	457607	9.08	408098	11.84
UPPER LIMIT	525000	8.7	915214	9.582	816196	12.341
LOWER LIMIT	131250	7.7	228804	8.582	204049	11.341
EPA SAMPLE NO.						
TW-1125-1	150985	8.21	353552	9.08	347429	11.84
TW-1125-2	203666	8.21	472280	9.08	460828	11.84
TB	177560	8.21	408212	9.08	395290	11.84
TB	168109	8.21	386817	9.08	372510	11.84
VN1201WBL01	188076	8.20	435319	9.08	418422	11.84
VN1201WBS01	256678	8.21	487472	9.08	435019	11.84
VN1201WBSD01	257123	8.21	472668	9.08	413438	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:	<u>Alliance</u>	Contract:	<u>REMI01</u>
Lab Code:	<u>ACE</u>	SDG NO.:	<u>Q3743</u>
Lab File ID:	<u>VN088370.D</u>	Date Analyzed:	<u>12/01/2025</u>
Instrument ID:	<u>MSVOA_N</u>	Time Analyzed:	<u>09:58</u>
GC Column:	<u>RXI-624</u>	ID:	<u>0.25</u> (mm)
		Heated Purge:	(Y/N) <u>N</u>

	IS4 AREA #	RT #				
12 HOUR STD	197460	13.765				
UPPER LIMIT	394920	14.265				
LOWER LIMIT	98730	13.265				
EPA SAMPLE NO.						
TW-1125-1	164080	13.77				
TW-1125-2	215787	13.77				
TB	183327	13.77				
TB	175948	13.77				
VN1201WBL01	198799	13.77				
VN1201WBS01	207303	13.77				
VN1201WBSD01	196888	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.



SAMPLE DATA

Report of Analysis

Client: Remington & Vernick
Project: Saddler Property
Client Sample ID: TW-1125-1
Lab Sample ID: Q3743-01
Analytical Method: 8260D
Sample Wt/Vol: 5 mL

Level: LOW
Final Vol: 5000 uL

Date Collected: 11/26/25
Date Received: 11/26/25
SDG No.: Q3743
Matrix: Water
% Solid: 0
Test: VOC-TCLVOA-10

CAS Number	Parameter	Conc.	Qua.	DF	MDL	LOQ / CRQL	Units	Date Ana.	BatchID
TARGETS									
75-71-8	Dichlorodifluoromethane	0.22	U	1	0.22	1.00	ug/L	12/01/25 18:20	VN120125
74-87-3	Chloromethane	0.32	U	1	0.32	1.00	ug/L	12/01/25 18:20	VN120125
75-01-4	Vinyl Chloride	0.26	U	1	0.26	1.00	ug/L	12/01/25 18:20	VN120125
74-83-9	Bromomethane	1.40	U	1	1.40	5.00	ug/L	12/01/25 18:20	VN120125
75-00-3	Chloroethane	0.47	U	1	0.47	1.00	ug/L	12/01/25 18:20	VN120125
75-69-4	Trichlorofluoromethane	0.33	U	1	0.33	1.00	ug/L	12/01/25 18:20	VN120125
76-13-1	1,1,2-Trichlorotrifluoroethane	0.25	U	1	0.25	1.00	ug/L	12/01/25 18:20	VN120125
75-35-4	1,1-Dichloroethene	0.23	U	1	0.23	1.00	ug/L	12/01/25 18:20	VN120125
67-64-1	Acetone	21.6		1	1.50	5.00	ug/L	12/01/25 18:20	VN120125
75-15-0	Carbon Disulfide	0.21	U	1	0.21	1.00	ug/L	12/01/25 18:20	VN120125
1634-04-4	Methyl tert-butyl Ether	0.16	U	1	0.16	1.00	ug/L	12/01/25 18:20	VN120125
79-20-9	Methyl Acetate	0.27	U	1	0.27	1.00	ug/L	12/01/25 18:20	VN120125
75-09-2	Methylene Chloride	0.28	U	1	0.28	1.00	ug/L	12/01/25 18:20	VN120125
156-60-5	trans-1,2-Dichloroethene	0.23	U	1	0.23	1.00	ug/L	12/01/25 18:20	VN120125
75-34-3	1,1-Dichloroethane	0.23	U	1	0.23	1.00	ug/L	12/01/25 18:20	VN120125
110-82-7	Cyclohexane	1.50	U	1	1.50	5.00	ug/L	12/01/25 18:20	VN120125
78-93-3	2-Butanone	0.98	U	1	0.98	5.00	ug/L	12/01/25 18:20	VN120125
56-23-5	Carbon Tetrachloride	0.25	U	1	0.25	1.00	ug/L	12/01/25 18:20	VN120125
156-59-2	cis-1,2-Dichloroethene	0.19	U	1	0.19	1.00	ug/L	12/01/25 18:20	VN120125
74-97-5	Bromochloromethane	0.22	U	1	0.22	1.00	ug/L	12/01/25 18:20	VN120125
67-66-3	Chloroform	0.25	U	1	0.25	1.00	ug/L	12/01/25 18:20	VN120125
71-55-6	1,1,1-Trichloroethane	0.20	U	1	0.20	1.00	ug/L	12/01/25 18:20	VN120125
108-87-2	Methylcyclohexane	0.16	U	1	0.16	1.00	ug/L	12/01/25 18:20	VN120125
71-43-2	Benzene	0.15	U	1	0.15	1.00	ug/L	12/01/25 18:20	VN120125
107-06-2	1,2-Dichloroethane	0.22	U	1	0.22	1.00	ug/L	12/01/25 18:20	VN120125
79-01-6	Trichloroethene	0.090	U	1	0.090	1.00	ug/L	12/01/25 18:20	VN120125
78-87-5	1,2-Dichloropropane	0.20	U	1	0.20	1.00	ug/L	12/01/25 18:20	VN120125
75-27-4	Bromodichloromethane	0.22	U	1	0.22	1.00	ug/L	12/01/25 18:20	VN120125
108-10-1	4-Methyl-2-Pentanone	0.68	U	1	0.68	5.00	ug/L	12/01/25 18:20	VN120125
108-88-3	Toluene	0.14	U	1	0.14	1.00	ug/L	12/01/25 18:20	VN120125
10061-02-6	t-1,3-Dichloropropene	0.17	U	1	0.17	1.00	ug/L	12/01/25 18:20	VN120125
10061-01-5	cis-1,3-Dichloropropene	0.16	U	1	0.16	1.00	ug/L	12/01/25 18:20	VN120125
79-00-5	1,1,2-Trichloroethane	0.21	U	1	0.21	1.00	ug/L	12/01/25 18:20	VN120125
591-78-6	2-Hexanone	0.89	U	1	0.89	5.00	ug/L	12/01/25 18:20	VN120125
124-48-1	Dibromochloromethane	0.18	U	1	0.18	1.00	ug/L	12/01/25 18:20	VN120125
106-93-4	1,2-Dibromoethane	0.15	U	1	0.15	1.00	ug/L	12/01/25 18:20	VN120125
127-18-4	Tetrachloroethene	0.23	U	1	0.23	1.00	ug/L	12/01/25 18:20	VN120125
108-90-7	Chlorobenzene	0.12	U	1	0.12	1.00	ug/L	12/01/25 18:20	VN120125
100-41-4	Ethyl Benzene	0.13	U	1	0.13	1.00	ug/L	12/01/25 18:20	VN120125
179601-23-1	m/p-Xylenes	0.24	U	1	0.24	2.00	ug/L	12/01/25 18:20	VN120125
95-47-6	o-Xylene	0.12	U	1	0.12	1.00	ug/L	12/01/25 18:20	VN120125
100-42-5	Styrene	0.15	U	1	0.15	1.00	ug/L	12/01/25 18:20	VN120125
75-25-2	Bromoform	0.19	U	1	0.19	1.00	ug/L	12/01/25 18:20	VN120125



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

Report of Analysis

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Lab Sample ID: Q3743-01
Analytical Method: 8260D
Sample Wt/Vol: 5 mL

Level: LOW
Final Vol: 5000 uL

Date Collected: 11/26/25
Date Received: 11/26/25
SDG No.: Q3743
Matrix: Water
% Solid: 0
Test: VOC-TCLVOA-10

CAS Number	Parameter	Conc.	Qua.	DF	MDL	LOQ / CRQL	Units	Date Ana.	BatchID
98-82-8	Isopropylbenzene	0.12	U	1	0.12	1.00	ug/L	12/01/25 18:20	VN120125
79-34-5	1,1,2,2-Tetrachloroethane	0.26	U	1	0.26	1.00	ug/L	12/01/25 18:20	VN120125
541-73-1	1,3-Dichlorobenzene	0.16	U	1	0.16	1.00	ug/L	12/01/25 18:20	VN120125
106-46-7	1,4-Dichlorobenzene	0.19	U	1	0.19	1.00	ug/L	12/01/25 18:20	VN120125
95-50-1	1,2-Dichlorobenzene	0.16	U	1	0.16	1.00	ug/L	12/01/25 18:20	VN120125
96-12-8	1,2-Dibromo-3-Chloropropane	0.53	U	1	0.53	1.00	ug/L	12/01/25 18:20	VN120125
120-82-1	1,2,4-Trichlorobenzene	0.20	U	1	0.20	1.00	ug/L	12/01/25 18:20	VN120125
87-61-6	1,2,3-Trichlorobenzene	0.20	U	1	0.20	1.00	ug/L	12/01/25 18:20	VN120125

SURROGATES

17060-07-0	1,2-Dichloroethane-d4	61.7			74 - 125	123%	SPK: 50
1868-53-7	Dibromofluoromethane	53.8			75 - 124	108%	SPK: 50
2037-26-5	Toluene-d8	50.2			86 - 113	100%	SPK: 50
460-00-4	4-Bromofluorobenzene	53.4			77 - 121	107%	SPK: 50

INTERNAL STANDARDS

INTERNAL STANDARDS		Area Count
363-72-4	Pentafluorobenzene	151000
540-36-3	1,4-Difluorobenzene	354000
3114-55-4	Chlorobenzene-d5	347000
3855-82-1	1,4-Dichlorobenzene-d4	164000

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\VN120125\
 Data File : VN088392.D
 Acq On : 01 Dec 2025 18:20
 Operator : JC\MD
 Sample : Q3743-01
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 TW-1125-1

Quant Time: Dec 02 00:18:44 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)

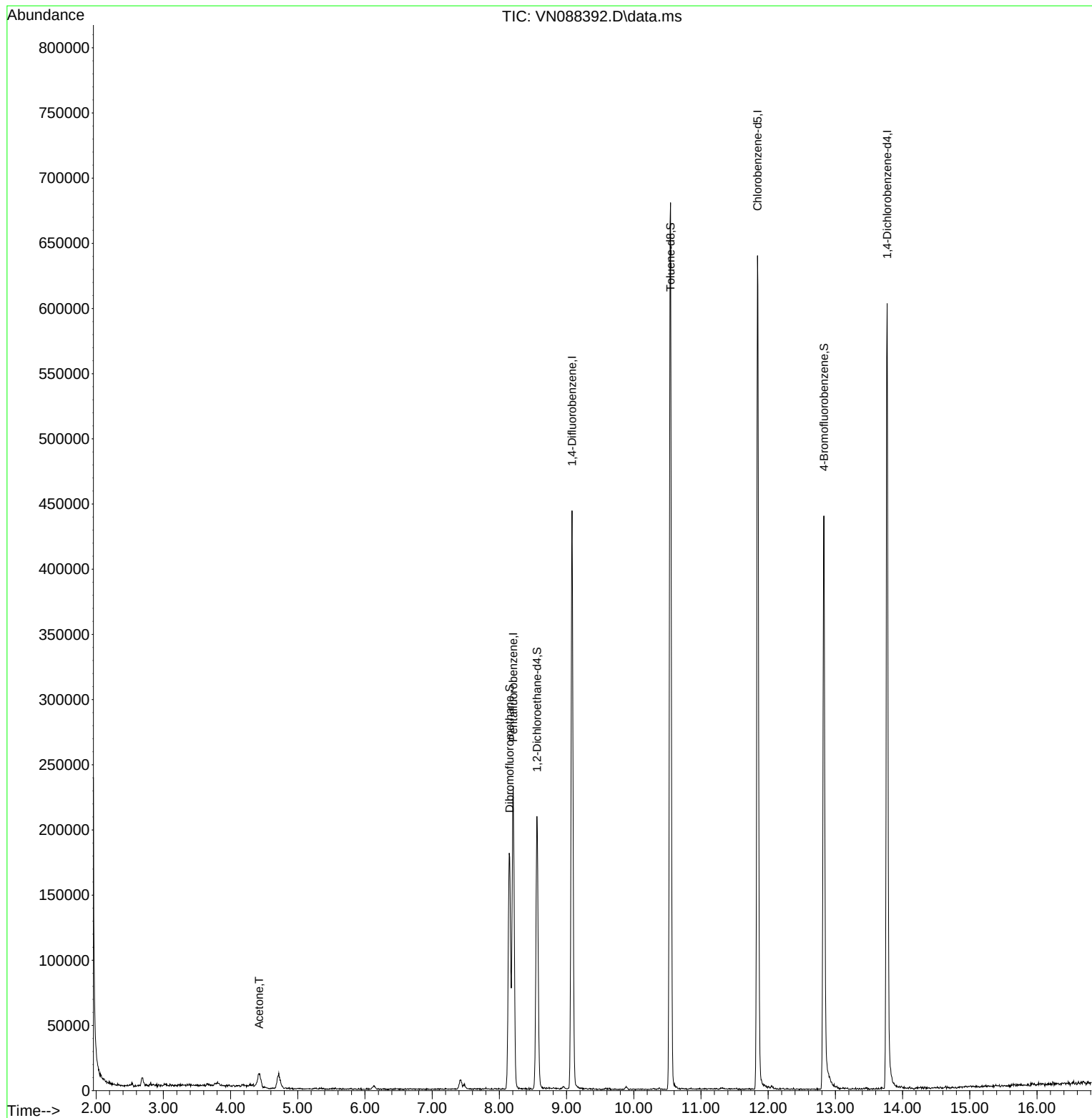
Internal Standards						
1) Pentafluorobenzene	8.206	168	150985	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.082	114	353552	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.841	117	347429	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.770	152	164080	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.559	65	175368	61.657	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	= 123.320%	
35) Dibromofluoromethane	8.153	113	131731	53.766	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 107.540%	
50) Toluene-d8	10.547	98	458002	50.240	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	= 100.480%	
62) 4-Bromofluorobenzene	12.829	95	166885	53.353	ug/l	0.00
Spiked Amount	50.000	Range	77 - 121	Recovery	= 106.700%	
Target Compounds						Qvalue
16) Acetone	4.424	43	22262	21.642	ug/l	97

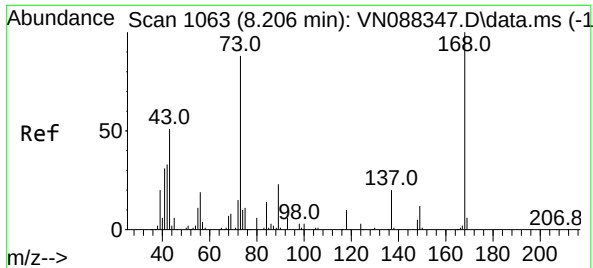
(#) = qualifier out of range (m) = manual integration (+) = signals summed

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN120125\
Data File : VN088392.D
Acq On : 01 Dec 2025 18:20
Operator : JC\MD
Sample : Q3743-01
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 25 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
TW-1125-1

Quant Time: Dec 02 00:18:44 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

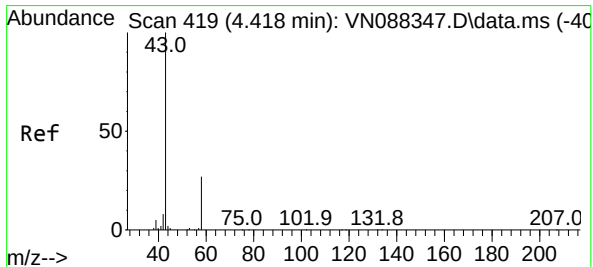
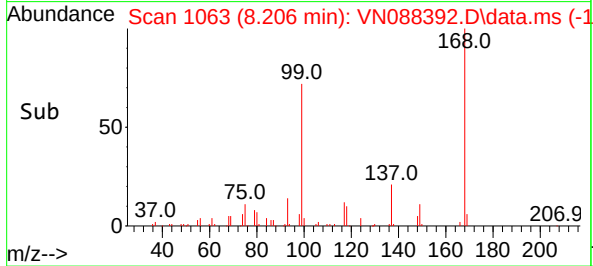
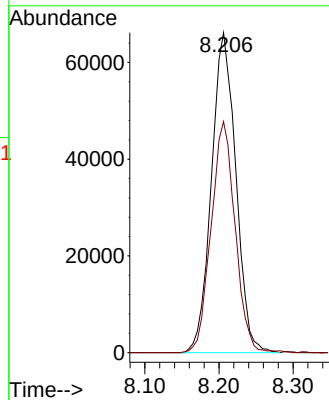
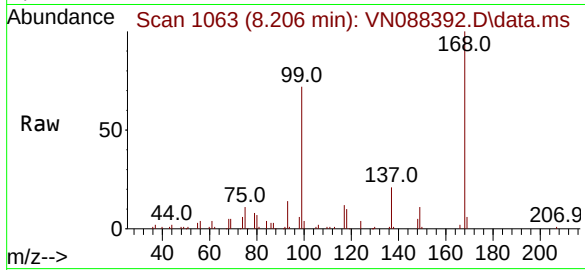




#1
Pentafluorobenzene
Concen: 50.000 ug/l
RT: 8.206 min Scan# 1063
Delta R.T. 0.000 min
Lab File: VN088392.D
Acq: 01 Dec 2025 18:20

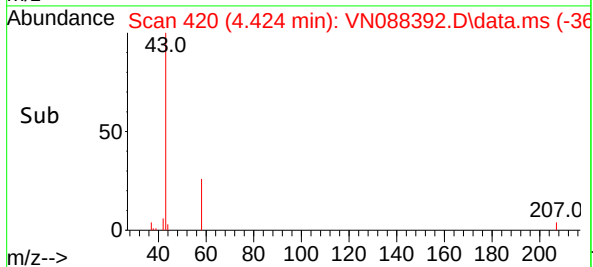
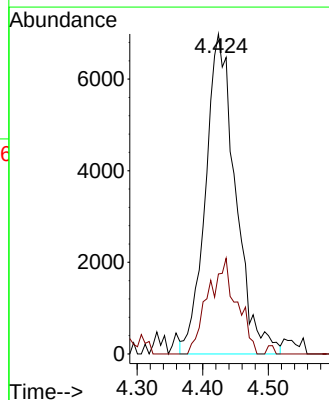
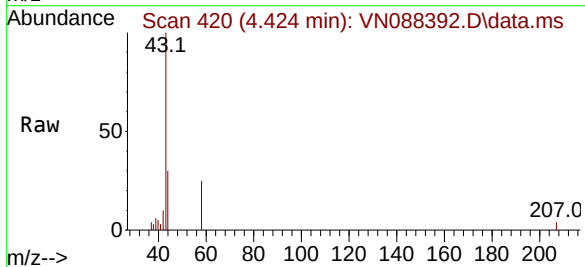
Instrument :
MSVOA_N
ClientSampleId :
TW-1125-1

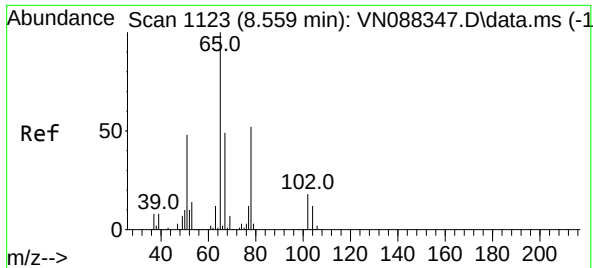
Tgt Ion:168 Resp: 150985
Ion Ratio Lower Upper
168 100
99 72.2 57.1 85.7



#16
Acetone
Concen: 21.642 ug/l
RT: 4.424 min Scan# 420
Delta R.T. 0.006 min
Lab File: VN088392.D
Acq: 01 Dec 2025 18:20

Tgt Ion: 43 Resp: 22262
Ion Ratio Lower Upper
43 100
58 25.6 21.7 32.5





#33

1,2-Dichloroethane-d4

Concen: 61.657 ug/l

RT: 8.559 min Scan# 1123

Delta R.T. 0.000 min

Lab File: VN088392.D

Acq: 01 Dec 2025 18:20

Instrument :

MSVOA_N

ClientSampleId :

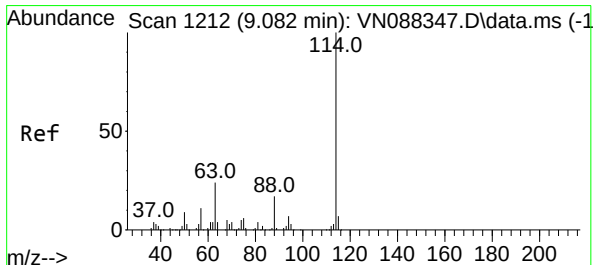
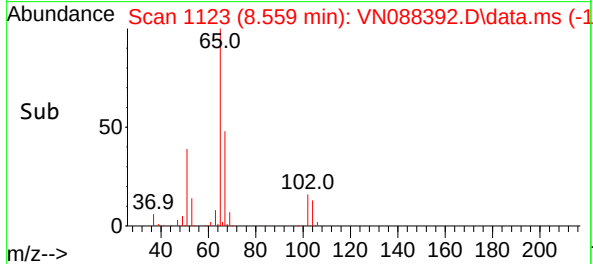
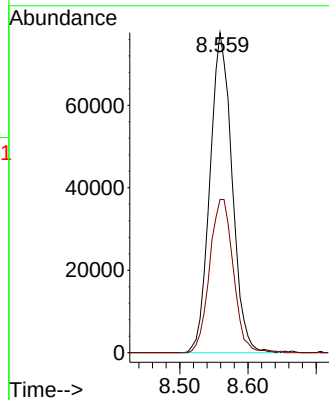
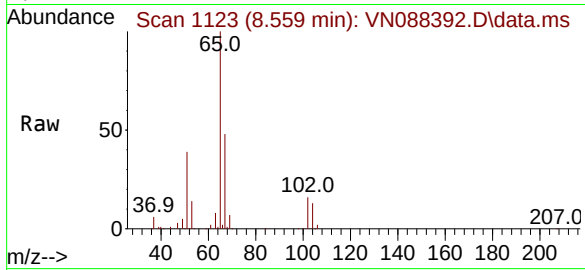
TW-1125-1

Tgt Ion: 65 Resp: 175368

Ion Ratio Lower Upper

65 100

67 51.1 0.0 100.8



#34

1,4-Difluorobenzene

Concen: 50.000 ug/l

RT: 9.082 min Scan# 1212

Delta R.T. 0.000 min

Lab File: VN088392.D

Acq: 01 Dec 2025 18:20

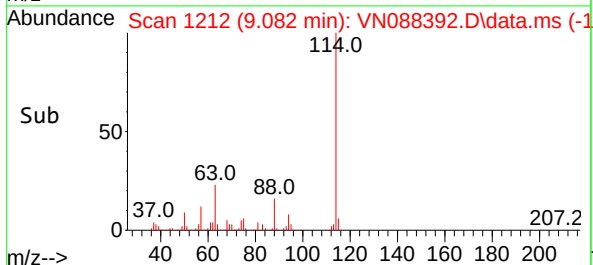
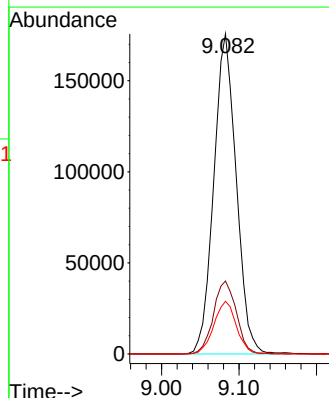
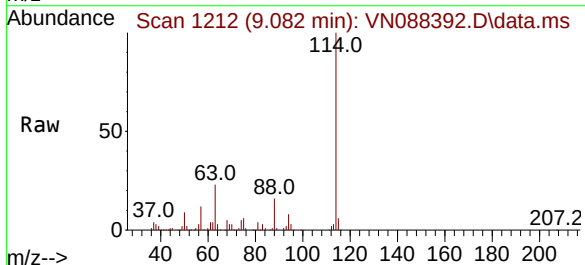
Tgt Ion: 114 Resp: 353552

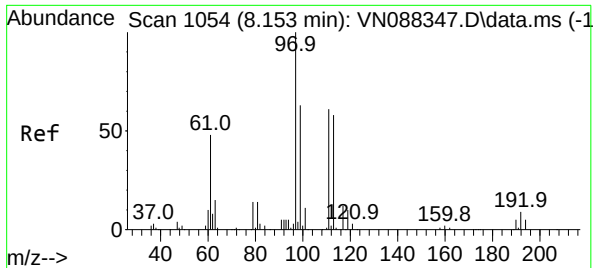
Ion Ratio Lower Upper

114 100

63 22.8 0.0 49.0

88 16.4 0.0 34.0





#35

Dibromofluoromethane

Concen: 53.766 ug/l

RT: 8.153 min Scan# 1054

Delta R.T. 0.000 min

Lab File: VN088392.D

Acq: 01 Dec 2025 18:20

Instrument :

MSVOA_N

ClientSampleId :

TW-1125-1

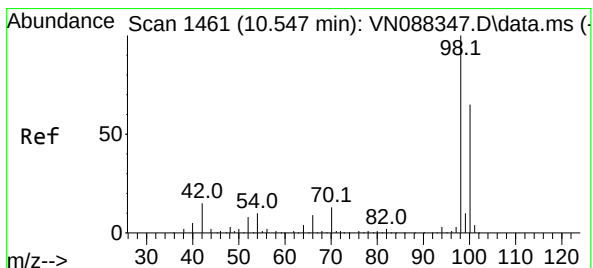
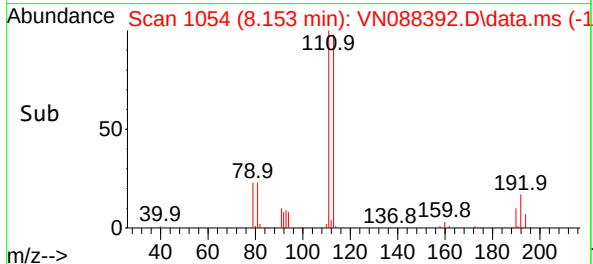
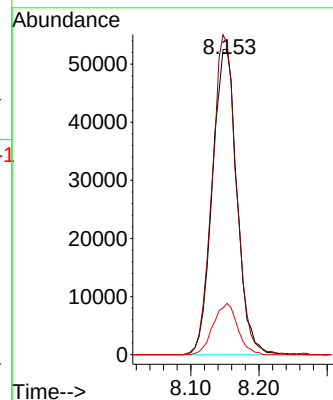
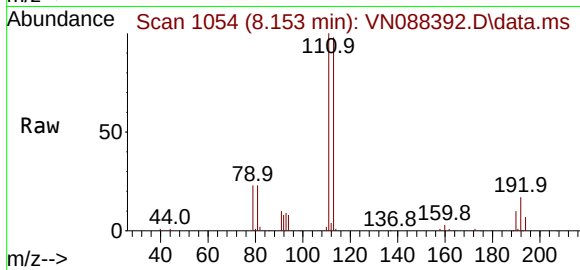
Tgt Ion:113 Resp: 131731

Ion Ratio Lower Upper

113 100

111 102.9 82.5 123.7

192 16.7 12.8 19.2



#50

Toluene-d8

Concen: 50.240 ug/l

RT: 10.547 min Scan# 1461

Delta R.T. 0.000 min

Lab File: VN088392.D

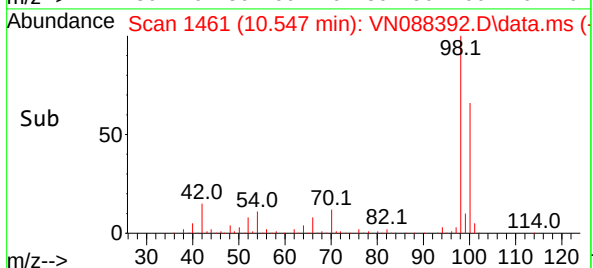
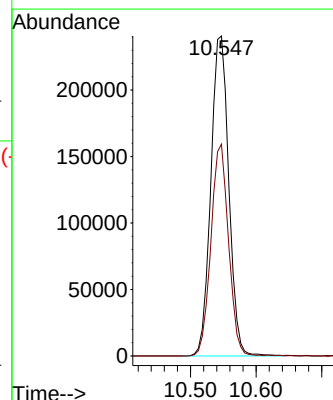
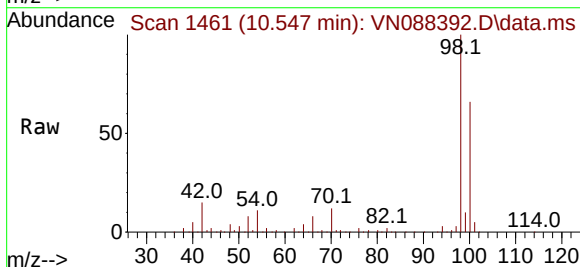
Acq: 01 Dec 2025 18:20

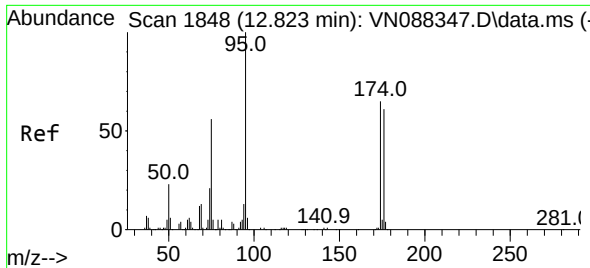
Tgt Ion: 98 Resp: 458002

Ion Ratio Lower Upper

98 100

100 64.8 52.2 78.2

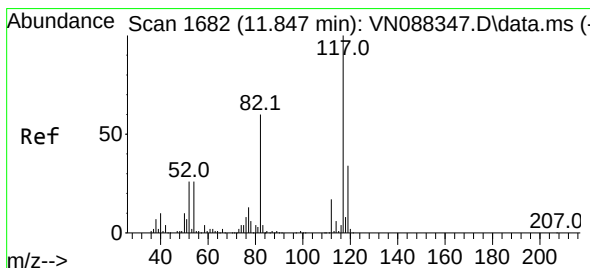
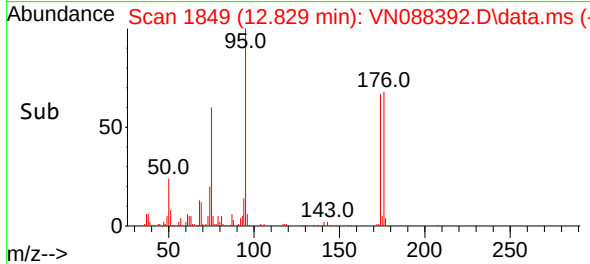
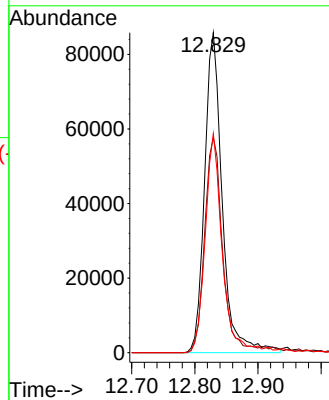
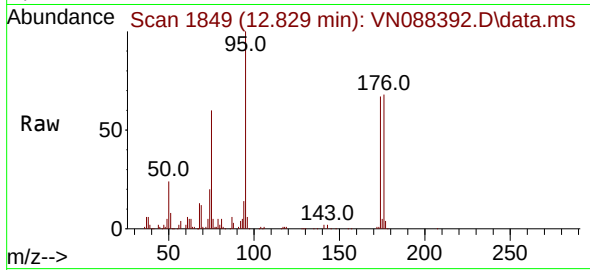




#62
4-Bromofluorobenzene
Concen: 53.353 ug/l
RT: 12.829 min Scan# 1848
Delta R.T. 0.006 min
Lab File: VN088392.D
Acq: 01 Dec 2025 18:20

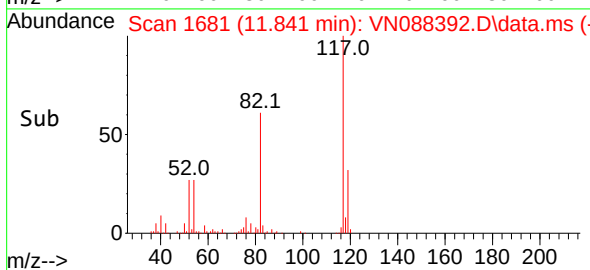
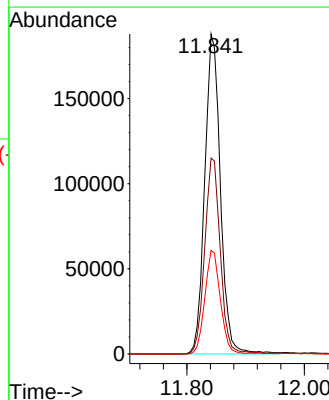
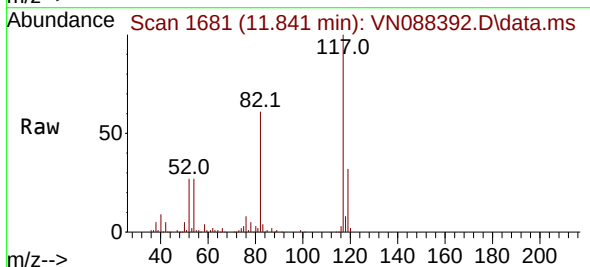
Instrument :
MSVOA_N
ClientSampleId :
TW-1125-1

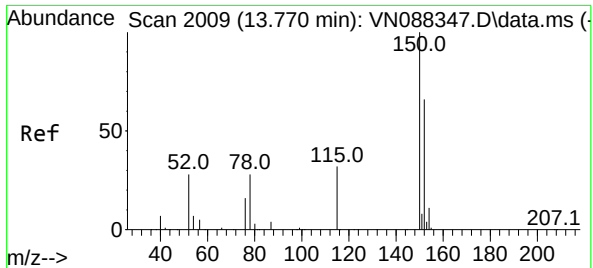
Tgt Ion: 95 Resp: 166885
Ion Ratio Lower Upper
95 100
174 67.7 0.0 139.0
176 66.4 0.0 132.4



#63
Chlorobenzene-d5
Concen: 50.000 ug/l
RT: 11.841 min Scan# 1681
Delta R.T. -0.006 min
Lab File: VN088392.D
Acq: 01 Dec 2025 18:20

Tgt Ion: 117 Resp: 347429
Ion Ratio Lower Upper
117 100
82 61.2 47.8 71.8
119 32.4 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: VN088392.D

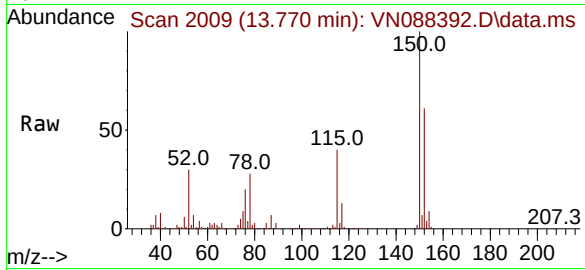
Acq: 01 Dec 2025 18:20

Instrument :

MSVOA_N

ClientSampleId :

TW-1125-1



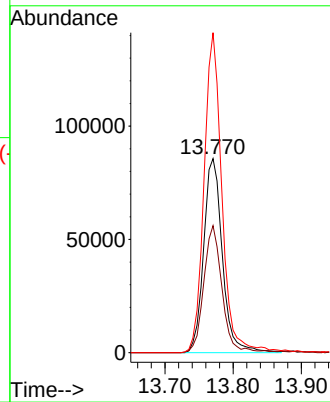
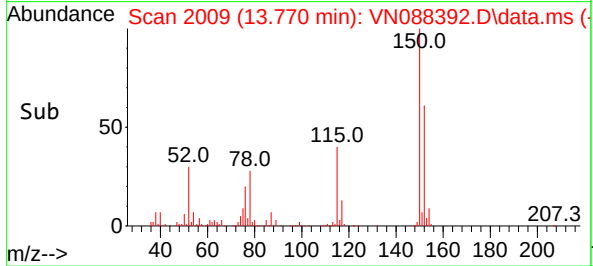
Tgt Ion:152 Resp: 164080

Ion Ratio Lower Upper

152 100

115 63.8 33.0 98.9

150 158.0 0.0 349.0



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN120125\
 Data File : VN088392.D
 Acq On : 01 Dec 2025 18:20
 Operator : JC\MD
 Sample : Q3743-01
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 TW-1125-1

Integration Parameters: RTEINT.P

Integrator: RTE

Smoothing : ON

Filtering: 5

Sampling : 1

Min Area: 3 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N112425W.M
 Title : SW846 8260

Signal : TIC: VN088392.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.683	120	124	132	rVB4	6295	14703	1.14%	0.213%
2	4.424	407	420	430	rVB3	10964	41128	3.20%	0.595%
3	4.718	460	470	483	rVB	11836	37976	2.96%	0.549%
4	7.424	921	930	935	rBV4	7022	18649	1.45%	0.270%
5	8.147	1042	1053	1058	rBV2	181041	453358	35.28%	6.556%
6	8.206	1058	1063	1077	rVB	232302	524149	40.79%	7.579%
7	8.559	1111	1123	1136	rBV2	209530	485377	37.78%	7.019%
8	9.082	1203	1212	1222	rBV	443172	905804	70.50%	13.098%
9	10.547	1452	1461	1470	rBV	680195	1284877	100.00%	18.580%
10	11.841	1673	1681	1694	rBV	639512	1180712	91.89%	17.073%
11	12.829	1841	1849	1862	rBV	439887	847908	65.99%	12.261%
12	13.770	2001	2009	2025	rBV	602142	1120846	87.23%	16.208%

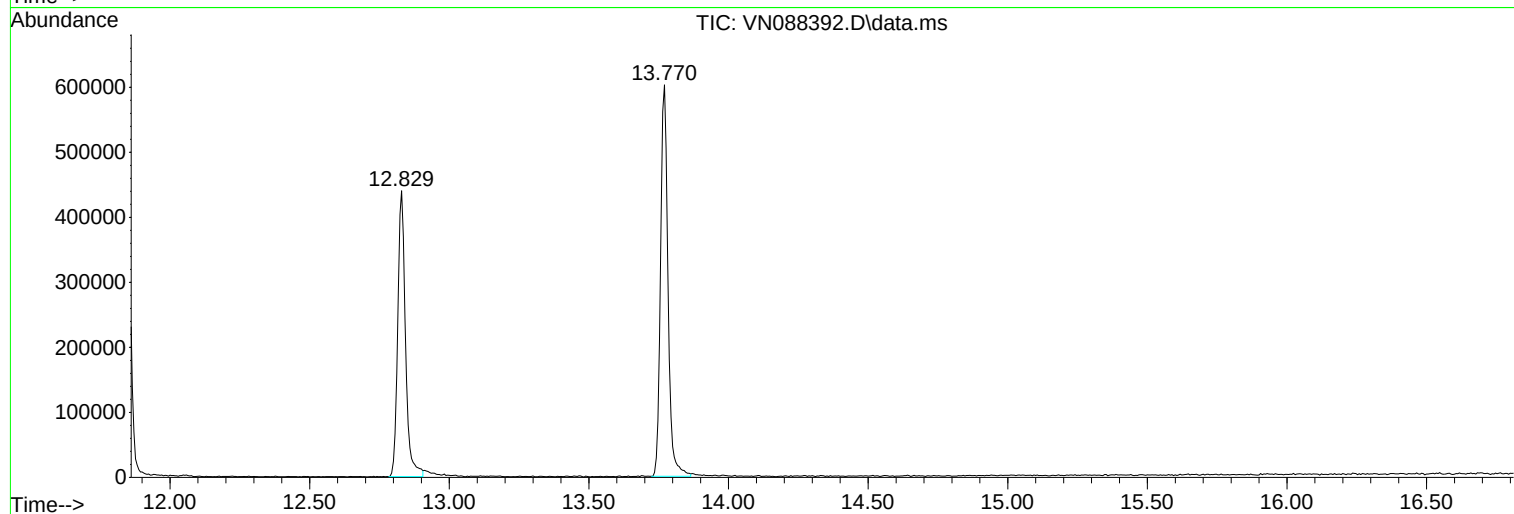
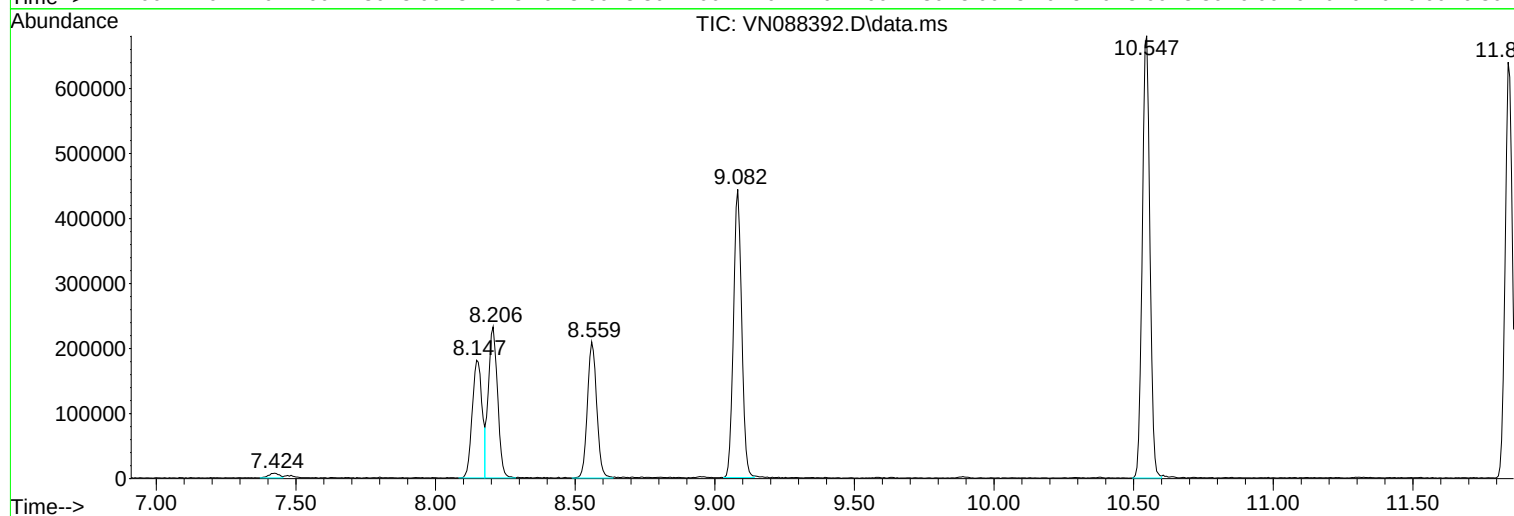
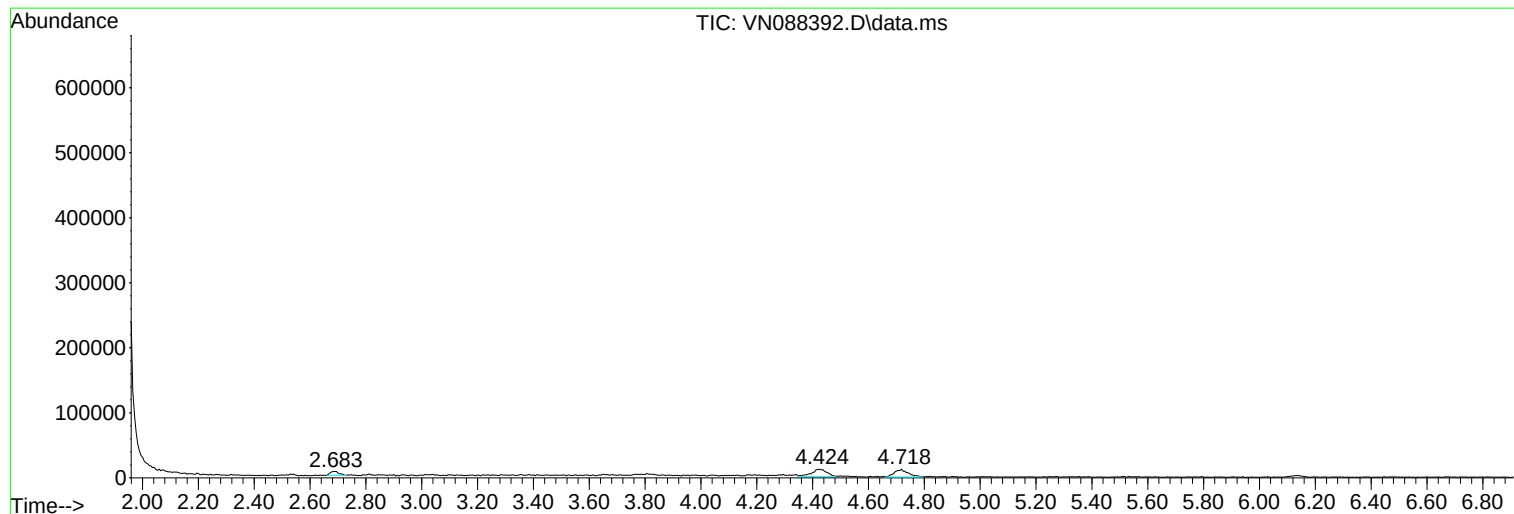
Sum of corrected areas: 6915487

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN120125\
Data File : VN088392.D
Acq On : 01 Dec 2025 18:20
Operator : JC\MD
Sample : Q3743-01
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 25 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
TW-1125-1

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N112425W.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN120125\
Data File : VN088392.D
Acq On : 01 Dec 2025 18:20
Operator : JC\MD
Sample : Q3743-01
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 25 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
TW-1125-1

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N112425W.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

No Library Search Compounds Detected

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN120125\
Data File : VN088392.D
Acq On : 01 Dec 2025 18:20
Operator : JC\MD
Sample : Q3743-01
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 25 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
TW-1125-1

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N112425W.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
------------------	----	---------	-------	----------	---	----	------	------

Report of Analysis

Client: Remington & Vernick
Project: Saddler Property
Client Sample ID: TW-1125-2
Lab Sample ID: Q3743-02
Analytical Method: 8260D
Sample Wt/Vol: 5 mL

Level: LOW
Final Vol: 5000 uL

Date Collected: 11/26/25
Date Received: 11/26/25
SDG No.: Q3743
Matrix: Water
% Solid: 0
Test: VOC-TCLVOA-10

CAS Number	Parameter	Conc.	Qua.	DF	MDL	LOQ / CRQL	Units	Date Ana.	BatchID
TARGETS									
75-71-8	Dichlorodifluoromethane	0.22	U	1	0.22	1.00	ug/L	12/01/25 18:41	VN120125
74-87-3	Chloromethane	0.32	U	1	0.32	1.00	ug/L	12/01/25 18:41	VN120125
75-01-4	Vinyl Chloride	0.26	U	1	0.26	1.00	ug/L	12/01/25 18:41	VN120125
74-83-9	Bromomethane	1.40	U	1	1.40	5.00	ug/L	12/01/25 18:41	VN120125
75-00-3	Chloroethane	0.47	U	1	0.47	1.00	ug/L	12/01/25 18:41	VN120125
75-69-4	Trichlorofluoromethane	0.33	U	1	0.33	1.00	ug/L	12/01/25 18:41	VN120125
76-13-1	1,1,2-Trichlorotrifluoroethane	0.25	U	1	0.25	1.00	ug/L	12/01/25 18:41	VN120125
75-35-4	1,1-Dichloroethene	0.23	U	1	0.23	1.00	ug/L	12/01/25 18:41	VN120125
67-64-1	Acetone	19.0		1	1.50	5.00	ug/L	12/01/25 18:41	VN120125
75-15-0	Carbon Disulfide	0.21	U	1	0.21	1.00	ug/L	12/01/25 18:41	VN120125
1634-04-4	Methyl tert-butyl Ether	0.16	U	1	0.16	1.00	ug/L	12/01/25 18:41	VN120125
79-20-9	Methyl Acetate	0.27	U	1	0.27	1.00	ug/L	12/01/25 18:41	VN120125
75-09-2	Methylene Chloride	0.28	U	1	0.28	1.00	ug/L	12/01/25 18:41	VN120125
156-60-5	trans-1,2-Dichloroethene	0.23	U	1	0.23	1.00	ug/L	12/01/25 18:41	VN120125
75-34-3	1,1-Dichloroethane	0.23	U	1	0.23	1.00	ug/L	12/01/25 18:41	VN120125
110-82-7	Cyclohexane	1.50	U	1	1.50	5.00	ug/L	12/01/25 18:41	VN120125
78-93-3	2-Butanone	0.98	U	1	0.98	5.00	ug/L	12/01/25 18:41	VN120125
56-23-5	Carbon Tetrachloride	0.25	U	1	0.25	1.00	ug/L	12/01/25 18:41	VN120125
156-59-2	cis-1,2-Dichloroethene	0.19	U	1	0.19	1.00	ug/L	12/01/25 18:41	VN120125
74-97-5	Bromochloromethane	0.22	U	1	0.22	1.00	ug/L	12/01/25 18:41	VN120125
67-66-3	Chloroform	0.25	U	1	0.25	1.00	ug/L	12/01/25 18:41	VN120125
71-55-6	1,1,1-Trichloroethane	0.20	U	1	0.20	1.00	ug/L	12/01/25 18:41	VN120125
108-87-2	Methylcyclohexane	0.16	U	1	0.16	1.00	ug/L	12/01/25 18:41	VN120125
71-43-2	Benzene	0.15	U	1	0.15	1.00	ug/L	12/01/25 18:41	VN120125
107-06-2	1,2-Dichloroethane	0.22	U	1	0.22	1.00	ug/L	12/01/25 18:41	VN120125
79-01-6	Trichloroethene	0.090	U	1	0.090	1.00	ug/L	12/01/25 18:41	VN120125
78-87-5	1,2-Dichloropropane	0.20	U	1	0.20	1.00	ug/L	12/01/25 18:41	VN120125
75-27-4	Bromodichloromethane	0.22	U	1	0.22	1.00	ug/L	12/01/25 18:41	VN120125
108-10-1	4-Methyl-2-Pentanone	0.68	U	1	0.68	5.00	ug/L	12/01/25 18:41	VN120125
108-88-3	Toluene	0.14	U	1	0.14	1.00	ug/L	12/01/25 18:41	VN120125
10061-02-6	t-1,3-Dichloropropene	0.17	U	1	0.17	1.00	ug/L	12/01/25 18:41	VN120125
10061-01-5	cis-1,3-Dichloropropene	0.16	U	1	0.16	1.00	ug/L	12/01/25 18:41	VN120125
79-00-5	1,1,2-Trichloroethane	0.21	U	1	0.21	1.00	ug/L	12/01/25 18:41	VN120125
591-78-6	2-Hexanone	0.89	U	1	0.89	5.00	ug/L	12/01/25 18:41	VN120125
124-48-1	Dibromochloromethane	0.18	U	1	0.18	1.00	ug/L	12/01/25 18:41	VN120125
106-93-4	1,2-Dibromoethane	0.15	U	1	0.15	1.00	ug/L	12/01/25 18:41	VN120125
127-18-4	Tetrachloroethene	0.23	U	1	0.23	1.00	ug/L	12/01/25 18:41	VN120125
108-90-7	Chlorobenzene	0.12	U	1	0.12	1.00	ug/L	12/01/25 18:41	VN120125
100-41-4	Ethyl Benzene	0.13	U	1	0.13	1.00	ug/L	12/01/25 18:41	VN120125
179601-23-1	m/p-Xylenes	0.24	U	1	0.24	2.00	ug/L	12/01/25 18:41	VN120125
95-47-6	o-Xylene	0.12	U	1	0.12	1.00	ug/L	12/01/25 18:41	VN120125
100-42-5	Styrene	0.15	U	1	0.15	1.00	ug/L	12/01/25 18:41	VN120125
75-25-2	Bromoform	0.19	U	1	0.19	1.00	ug/L	12/01/25 18:41	VN120125

Report of Analysis

Client:	Remington & Vernick	Date Collected:	11/26/25
Project:	Saddler Property	Date Received:	11/26/25
Client Sample ID:	TW-1125-2	SDG No.:	Q3743
Lab Sample ID:	Q3743-02	Matrix:	Water
Analytical Method:	8260D	% Solid:	0
Sample Wt/Vol:	5 mL	Test:	VOC-TCLVOA-10
	Level: LOW		
	Final Vol: 5000 uL		

CAS Number	Parameter	Conc.	Qua.	DF	MDL	LOQ / CRQL	Units	Date Ana.	BatchID
98-82-8	Isopropylbenzene	0.12	U	1	0.12	1.00	ug/L	12/01/25 18:41	VN120125
79-34-5	1,1,2,2-Tetrachloroethane	0.26	U	1	0.26	1.00	ug/L	12/01/25 18:41	VN120125
541-73-1	1,3-Dichlorobenzene	0.16	U	1	0.16	1.00	ug/L	12/01/25 18:41	VN120125
106-46-7	1,4-Dichlorobenzene	0.19	U	1	0.19	1.00	ug/L	12/01/25 18:41	VN120125
95-50-1	1,2-Dichlorobenzene	0.16	U	1	0.16	1.00	ug/L	12/01/25 18:41	VN120125
96-12-8	1,2-Dibromo-3-Chloropropane	0.53	U	1	0.53	1.00	ug/L	12/01/25 18:41	VN120125
120-82-1	1,2,4-Trichlorobenzene	0.20	U	1	0.20	1.00	ug/L	12/01/25 18:41	VN120125
87-61-6	1,2,3-Trichlorobenzene	0.20	U	1	0.20	1.00	ug/L	12/01/25 18:41	VN120125
SURROGATES									
17060-07-0	1,2-Dichloroethane-d4	61.2			74 - 125	122%	SPK: 50		
1868-53-7	Dibromofluoromethane	53.0			75 - 124	106%	SPK: 50		
2037-26-5	Toluene-d8	50.1			86 - 113	100%	SPK: 50		
460-00-4	4-Bromofluorobenzene	53.5			77 - 121	107%	SPK: 50		
INTERNAL STANDARDS									
		Area Count							
363-72-4	Pentafluorobenzene	204000							
540-36-3	1,4-Difluorobenzene	472000							
3114-55-4	Chlorobenzene-d5	461000							
3855-82-1	1,4-Dichlorobenzene-d4	216000							

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\VN120125\
 Data File : VN088393.D
 Acq On : 01 Dec 2025 18:41
 Operator : JC\MD
 Sample : Q3743-02
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 TW-1125-2

Quant Time: Dec 02 00:19:06 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)

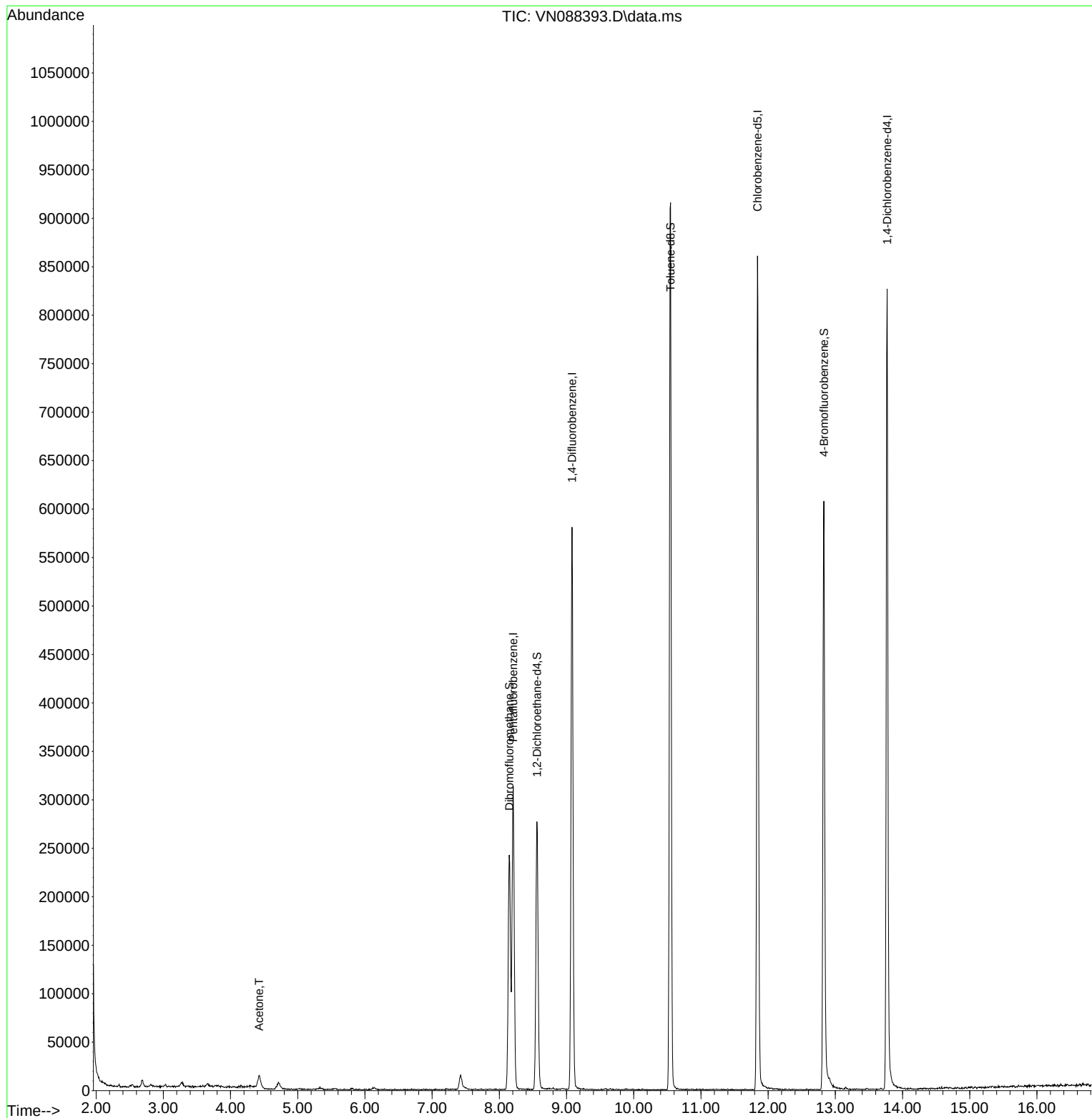
Internal Standards						
1) Pentafluorobenzene	8.206	168	203666	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.083	114	472280	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.841	117	460828	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.771	152	215787	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.559	65	234894	61.224	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	= 122.440%	
35) Dibromofluoromethane	8.147	113	173373	52.973	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 105.940%	
50) Toluene-d8	10.547	98	610228	50.110	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	= 100.220%	
62) 4-Bromofluorobenzene	12.829	95	223615	53.517	ug/l	0.00
Spiked Amount	50.000	Range	77 - 121	Recovery	= 107.040%	
Target Compounds						
16) Acetone	4.424	43	26336	18.980	ug/l	Qvalue 96

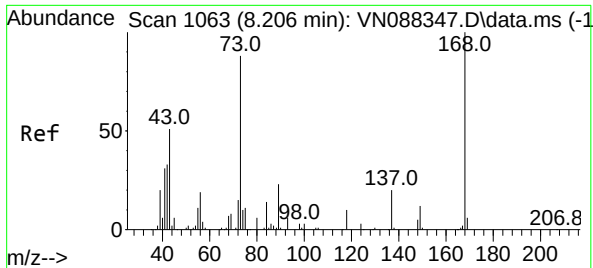
(#) = qualifier out of range (m) = manual integration (+) = signals summed

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN120125\
Data File : VN088393.D
Acq On : 01 Dec 2025 18:41
Operator : JC\MD
Sample : Q3743-02
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 26 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
TW-1125-2

Quant Time: Dec 02 00:19:06 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

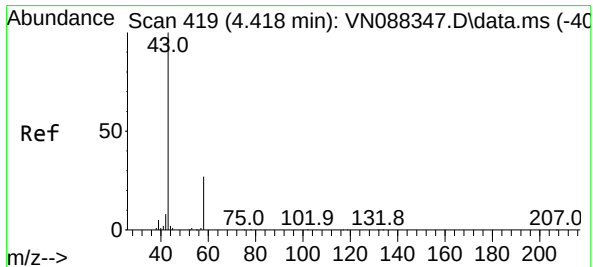
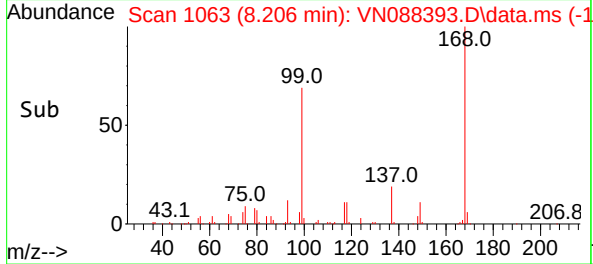
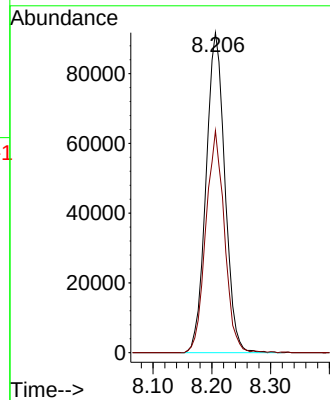
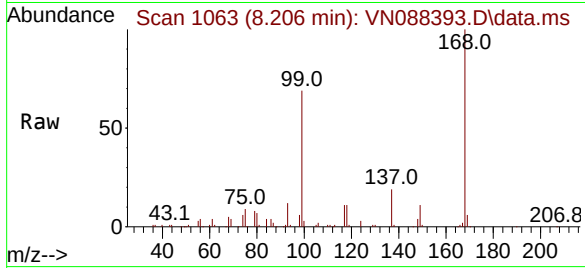




#1
Pentafluorobenzene
Concen: 50.000 ug/l
RT: 8.206 min Scan# 1063
Delta R.T. 0.000 min
Lab File: VN088393.D
Acq: 01 Dec 2025 18:41

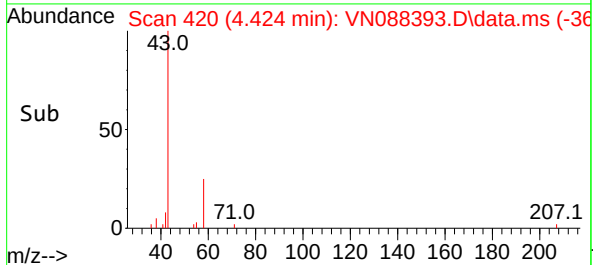
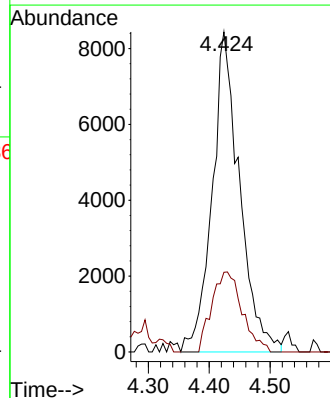
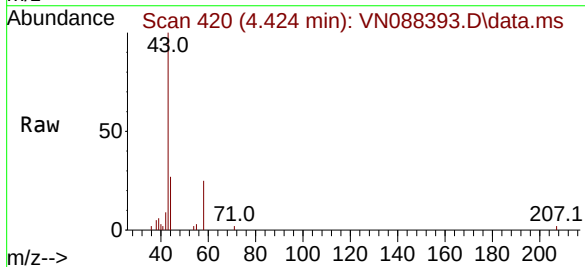
Instrument :
MSVOA_N
ClientSampleId :
TW-1125-2

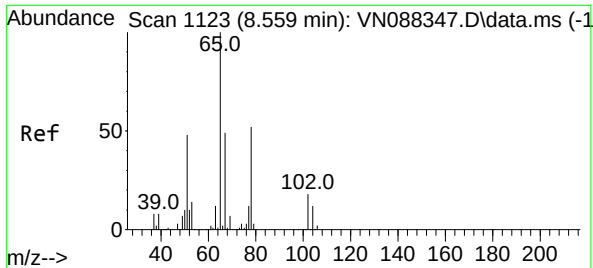
Tgt Ion:168 Resp: 203666
Ion Ratio Lower Upper
168 100
99 69.4 57.1 85.7



#16
Acetone
Concen: 18.980 ug/l
RT: 4.424 min Scan# 420
Delta R.T. 0.006 min
Lab File: VN088393.D
Acq: 01 Dec 2025 18:41

Tgt Ion: 43 Resp: 26336
Ion Ratio Lower Upper
43 100
58 24.9 21.7 32.5





#33

1,2-Dichloroethane-d4

Concen: 61.224 ug/l

RT: 8.559 min Scan# 1123

Delta R.T. 0.000 min

Lab File: VN088393.D

Acq: 01 Dec 2025 18:41

Instrument :

MSVOA_N

ClientSampleId :

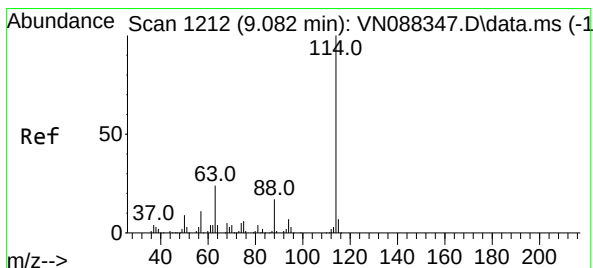
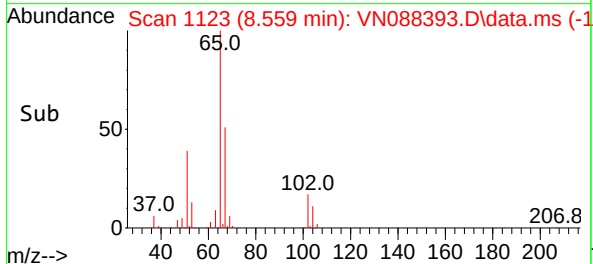
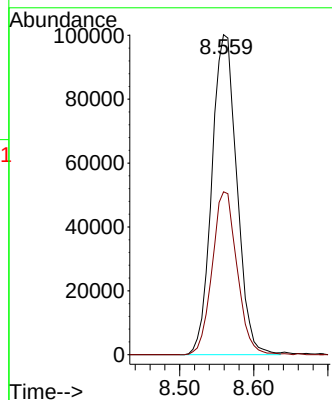
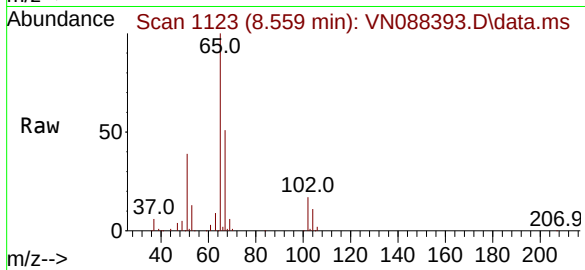
TW-1125-2

Tgt Ion: 65 Resp: 234894

Ion Ratio Lower Upper

65 100

67 50.6 0.0 100.8



#34

1,4-Difluorobenzene

Concen: 50.000 ug/l

RT: 9.083 min Scan# 1212

Delta R.T. 0.000 min

Lab File: VN088393.D

Acq: 01 Dec 2025 18:41

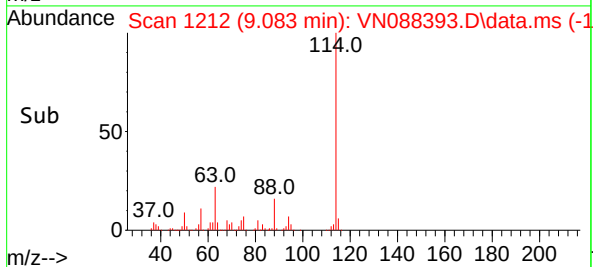
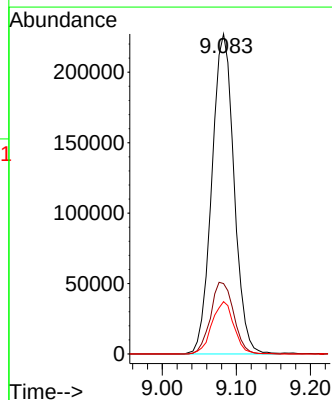
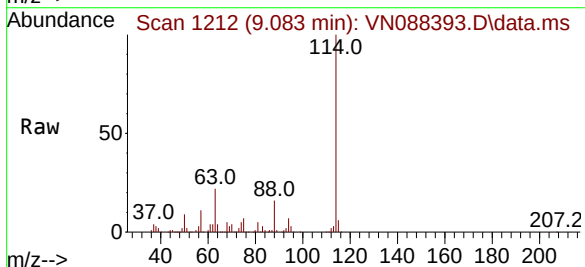
Tgt Ion: 114 Resp: 472280

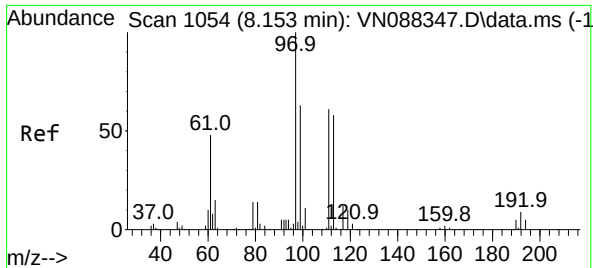
Ion Ratio Lower Upper

114 100

63 22.0 0.0 49.0

88 16.4 0.0 34.0





#35

Dibromofluoromethane

Concen: 52.973 ug/l

RT: 8.147 min Scan# 1053

Delta R.T. -0.006 min

Lab File: VN088393.D

Acq: 01 Dec 2025 18:41

Instrument :

MSVOA_N

ClientSampleId :

TW-1125-2

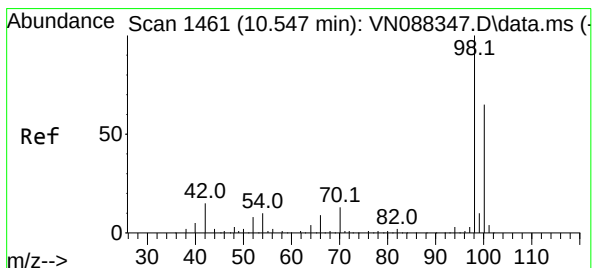
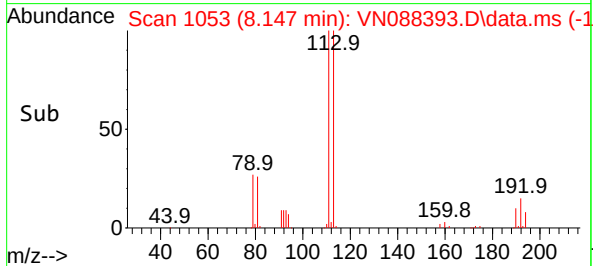
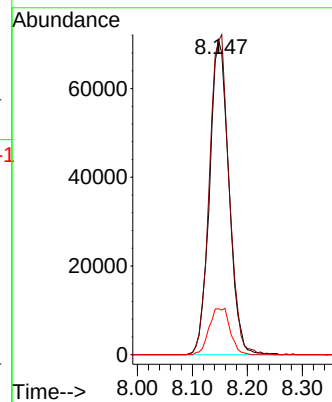
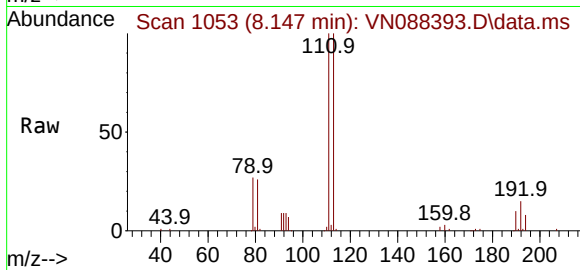
Tgt Ion:113 Resp: 173373

Ion Ratio Lower Upper

113 100

111 103.3 82.5 123.7

192 15.9 12.8 19.2



#50

Toluene-d8

Concen: 50.110 ug/l

RT: 10.547 min Scan# 1461

Delta R.T. 0.000 min

Lab File: VN088393.D

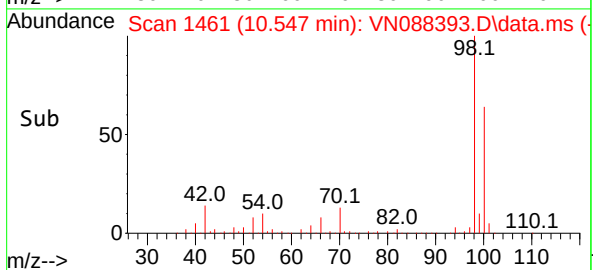
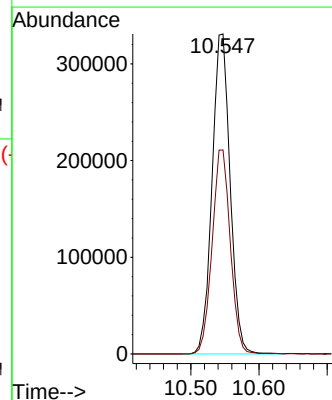
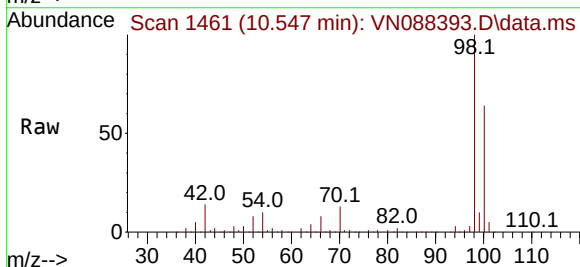
Acq: 01 Dec 2025 18:41

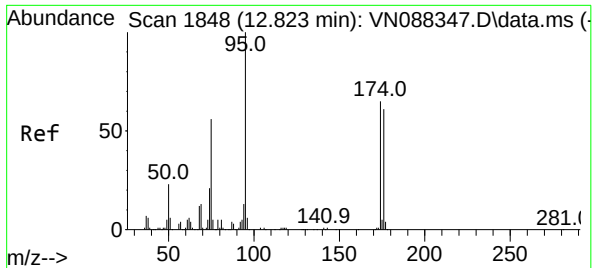
Tgt Ion: 98 Resp: 610228

Ion Ratio Lower Upper

98 100

100 64.3 52.2 78.2

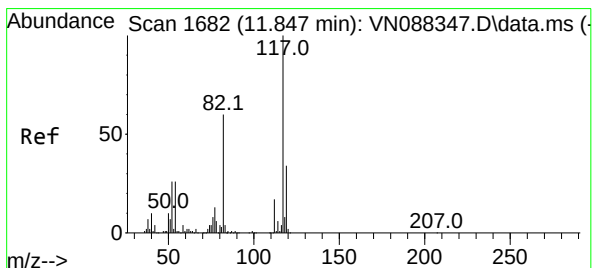
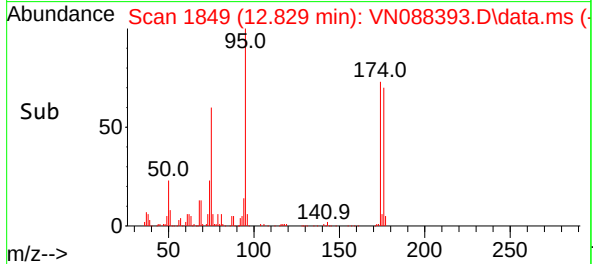
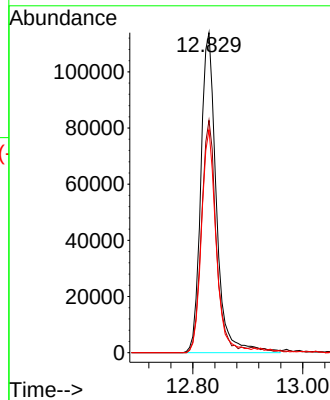
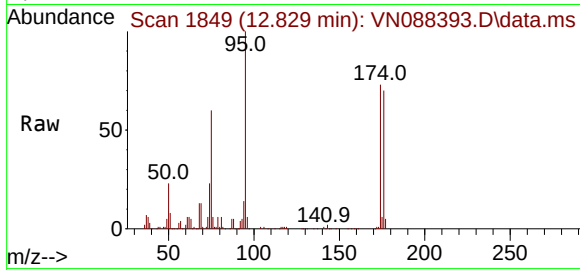




#62
4-Bromofluorobenzene
Concen: 53.517 ug/l
RT: 12.829 min Scan# 1848
Delta R.T. 0.006 min
Lab File: VN088393.D
Acq: 01 Dec 2025 18:41

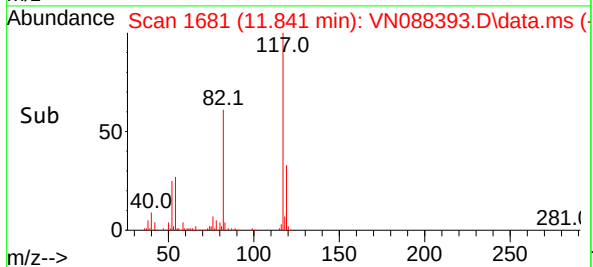
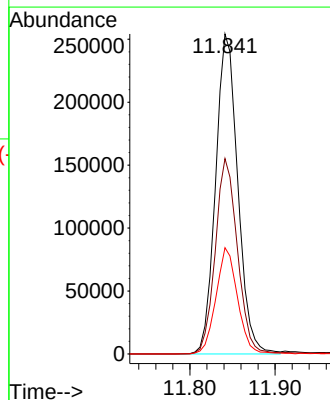
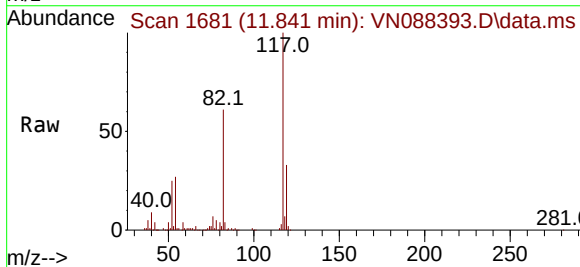
Instrument :
MSVOA_N
ClientSampleId :
TW-1125-2

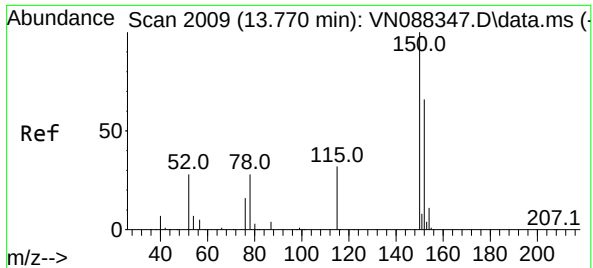
Tgt Ion: 95 Resp: 223615
Ion Ratio Lower Upper
95 100
174 69.2 0.0 139.0
176 66.0 0.0 132.4



#63
Chlorobenzene-d5
Concen: 50.000 ug/l
RT: 11.841 min Scan# 1681
Delta R.T. -0.006 min
Lab File: VN088393.D
Acq: 01 Dec 2025 18:41

Tgt Ion: 117 Resp: 460828
Ion Ratio Lower Upper
117 100
82 61.0 47.8 71.8
119 33.2 26.9 40.3





#72

1,4-Dichlorobenzene-d4

Concen: 50.000 ug/l

RT: 13.771 min Scan# 20

Delta R.T. 0.000 min

Lab File: VN088393.D

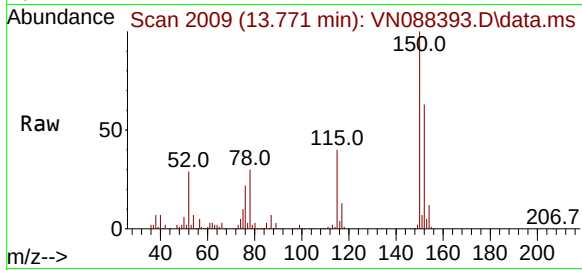
Acq: 01 Dec 2025 18:41

Instrument :

MSVOA_N

ClientSampleId :

TW-1125-2



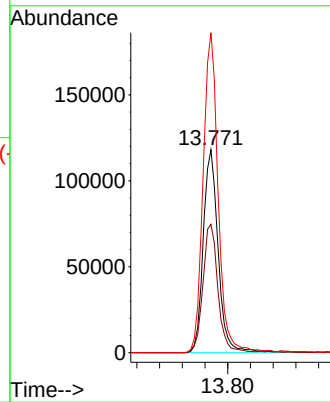
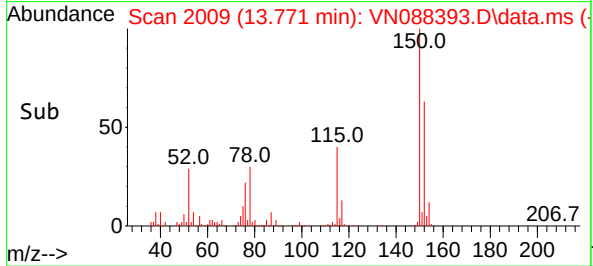
Tgt Ion:152 Resp: 215787

Ion Ratio Lower Upper

152 100

115 63.2 33.0 98.9

150 158.4 0.0 349.0



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN120125\
Data File : VN088393.D
Acq On : 01 Dec 2025 18:41
Operator : JC\MD
Sample : Q3743-02
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 26 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
TW-1125-2

Integration Parameters: RTEINT.P

Integrator: RTE

Smoothing : ON

Filtering: 5

Sampling : 1

Min Area: 3 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

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

Title : SW846 8260

Signal : TIC: VN088393.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.424	412	420	429	rVB2	12381	36002	2.10%	0.394%
2	7.424	921	930	942	rBV	14473	43216	2.53%	0.473%
3	8.147	1043	1053	1058	rBV2	241953	595850	34.82%	6.522%
4	8.206	1058	1063	1075	rVB	312188	686641	40.12%	7.515%
5	8.559	1112	1123	1135	rBV2	276489	640974	37.45%	7.016%
6	9.083	1203	1212	1225	rBV	579905	1212927	70.88%	13.276%
7	10.547	1452	1461	1472	rBV	915592	1711321	100.00%	18.731%
8	11.841	1673	1681	1695	rBV	860279	1575906	92.09%	17.249%
9	12.829	1839	1849	1861	rBV	607453	1144635	66.89%	12.528%
10	13.771	2001	2009	2023	rBV	825295	1488997	87.01%	16.297%

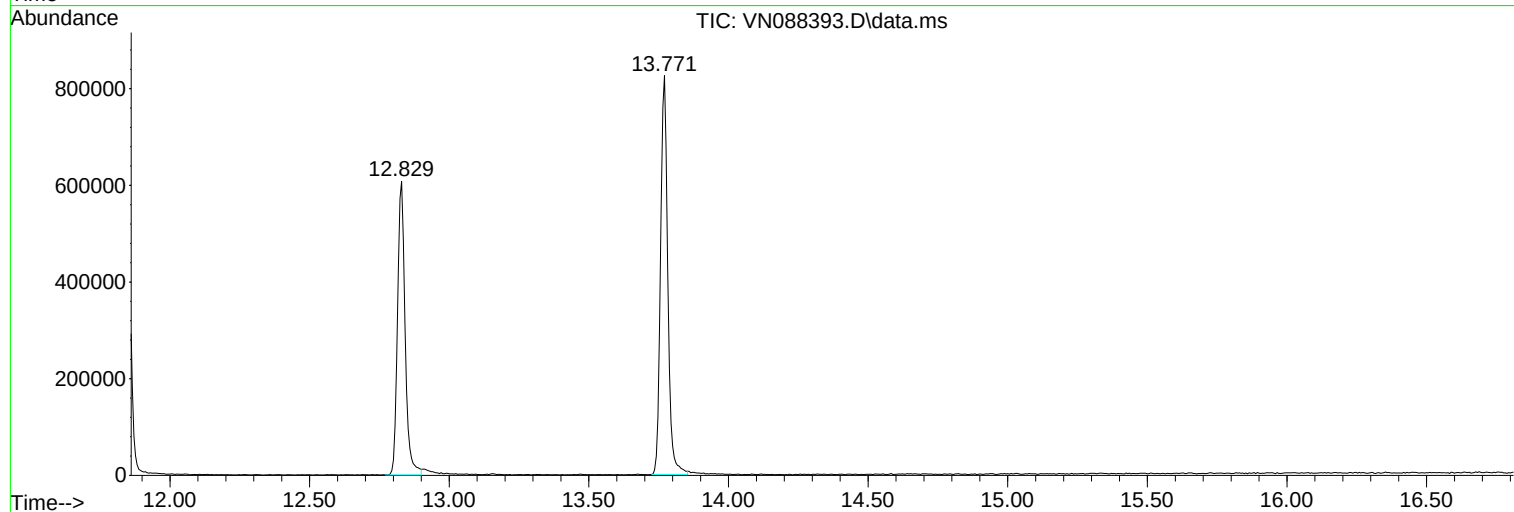
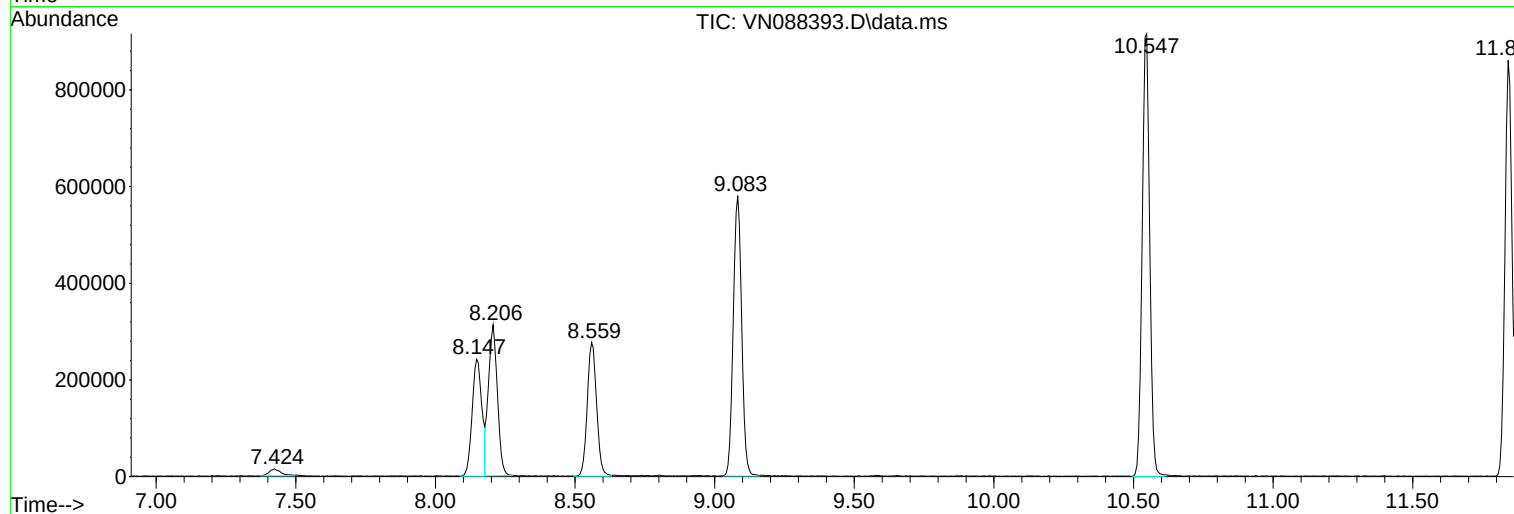
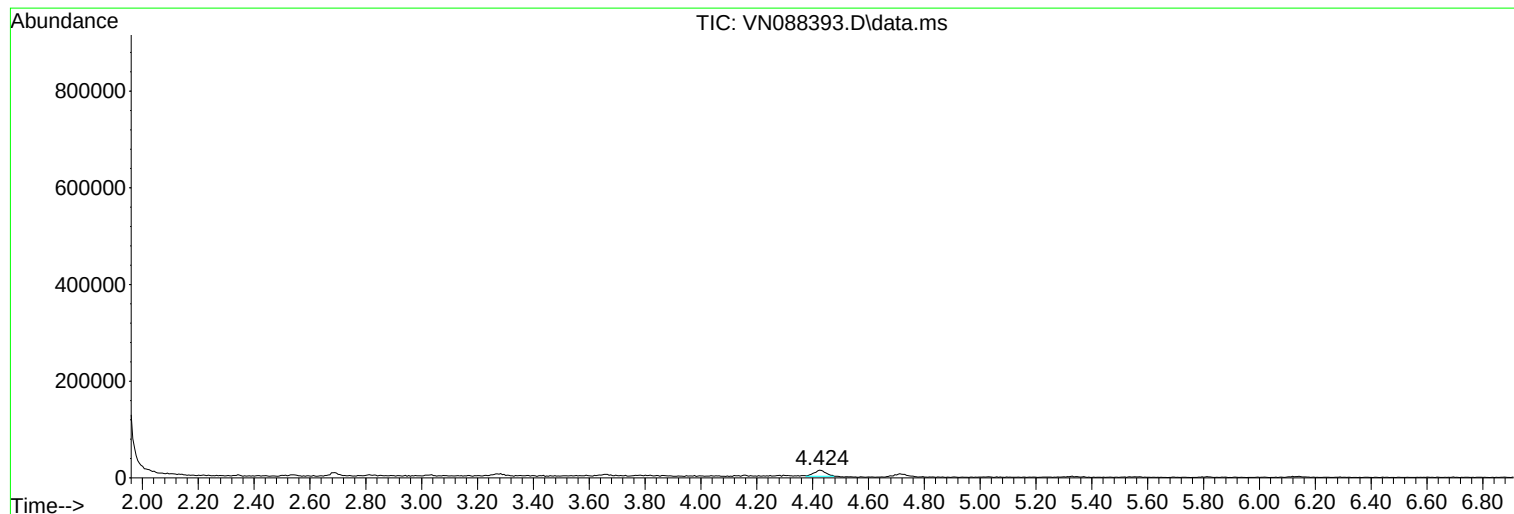
Sum of corrected areas: 9136469

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN120125\
Data File : VN088393.D
Acq On : 01 Dec 2025 18:41
Operator : JC\MD
Sample : Q3743-02
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 26 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
TW-1125-2

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N112425W.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN120125\
Data File : VN088393.D
Acq On : 01 Dec 2025 18:41
Operator : JC\MD
Sample : Q3743-02
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 26 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
TW-1125-2

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N112425W.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

No Library Search Compounds Detected

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN120125\
Data File : VN088393.D
Acq On : 01 Dec 2025 18:41
Operator : JC\MD
Sample : Q3743-02
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 26 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
TW-1125-2

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N112425W.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc

Report of Analysis

Client: Remington & Vernick
Project: Saddler Property
Client Sample ID: TB
Lab Sample ID: Q3743-03
Analytical Method: 8260D
Sample Wt/Vol: 5 mL

Level: LOW
Final Vol: 5000 uL

Date Collected: 11/26/25
Date Received: 11/26/25
SDG No.: Q3743
Matrix: Water
% Solid: 0
Test: VOC-TCLVOA-10

CAS Number	Parameter	Conc.	Qua.	DF	MDL	LOQ / CRQL	Units	Date Ana.	BatchID
TARGETS									
75-71-8	Dichlorodifluoromethane	0.22	U	1	0.22	1.00	ug/L	12/01/25 17:18	VN120125
74-87-3	Chloromethane	0.32	U	1	0.32	1.00	ug/L	12/01/25 17:18	VN120125
75-01-4	Vinyl Chloride	0.26	U	1	0.26	1.00	ug/L	12/01/25 17:18	VN120125
74-83-9	Bromomethane	1.40	U	1	1.40	5.00	ug/L	12/01/25 17:18	VN120125
75-00-3	Chloroethane	0.47	U	1	0.47	1.00	ug/L	12/01/25 17:18	VN120125
75-69-4	Trichlorofluoromethane	0.33	U	1	0.33	1.00	ug/L	12/01/25 17:18	VN120125
76-13-1	1,1,2-Trichlorotrifluoroethane	0.25	U	1	0.25	1.00	ug/L	12/01/25 17:18	VN120125
75-35-4	1,1-Dichloroethene	0.23	U	1	0.23	1.00	ug/L	12/01/25 17:18	VN120125
67-64-1	Acetone	1.50	U	1	1.50	5.00	ug/L	12/01/25 17:18	VN120125
75-15-0	Carbon Disulfide	0.21	U	1	0.21	1.00	ug/L	12/01/25 17:18	VN120125
1634-04-4	Methyl tert-butyl Ether	0.16	U	1	0.16	1.00	ug/L	12/01/25 17:18	VN120125
79-20-9	Methyl Acetate	0.27	U	1	0.27	1.00	ug/L	12/01/25 17:18	VN120125
75-09-2	Methylene Chloride	0.28	U	1	0.28	1.00	ug/L	12/01/25 17:18	VN120125
156-60-5	trans-1,2-Dichloroethene	0.23	U	1	0.23	1.00	ug/L	12/01/25 17:18	VN120125
75-34-3	1,1-Dichloroethane	0.23	U	1	0.23	1.00	ug/L	12/01/25 17:18	VN120125
110-82-7	Cyclohexane	1.50	U	1	1.50	5.00	ug/L	12/01/25 17:18	VN120125
78-93-3	2-Butanone	0.98	U	1	0.98	5.00	ug/L	12/01/25 17:18	VN120125
56-23-5	Carbon Tetrachloride	0.25	U	1	0.25	1.00	ug/L	12/01/25 17:18	VN120125
156-59-2	cis-1,2-Dichloroethene	0.19	U	1	0.19	1.00	ug/L	12/01/25 17:18	VN120125
74-97-5	Bromochloromethane	0.22	U	1	0.22	1.00	ug/L	12/01/25 17:18	VN120125
67-66-3	Chloroform	0.25	U	1	0.25	1.00	ug/L	12/01/25 17:18	VN120125
71-55-6	1,1,1-Trichloroethane	0.20	U	1	0.20	1.00	ug/L	12/01/25 17:18	VN120125
108-87-2	Methylcyclohexane	0.16	U	1	0.16	1.00	ug/L	12/01/25 17:18	VN120125
71-43-2	Benzene	0.15	U	1	0.15	1.00	ug/L	12/01/25 17:18	VN120125
107-06-2	1,2-Dichloroethane	0.22	U	1	0.22	1.00	ug/L	12/01/25 17:18	VN120125
79-01-6	Trichloroethene	0.090	U	1	0.090	1.00	ug/L	12/01/25 17:18	VN120125
78-87-5	1,2-Dichloropropane	0.20	U	1	0.20	1.00	ug/L	12/01/25 17:18	VN120125
75-27-4	Bromodichloromethane	0.22	U	1	0.22	1.00	ug/L	12/01/25 17:18	VN120125
108-10-1	4-Methyl-2-Pentanone	0.68	U	1	0.68	5.00	ug/L	12/01/25 17:18	VN120125
108-88-3	Toluene	0.14	U	1	0.14	1.00	ug/L	12/01/25 17:18	VN120125
10061-02-6	t-1,3-Dichloropropene	0.17	U	1	0.17	1.00	ug/L	12/01/25 17:18	VN120125
10061-01-5	cis-1,3-Dichloropropene	0.16	U	1	0.16	1.00	ug/L	12/01/25 17:18	VN120125
79-00-5	1,1,2-Trichloroethane	0.21	U	1	0.21	1.00	ug/L	12/01/25 17:18	VN120125
591-78-6	2-Hexanone	0.89	U	1	0.89	5.00	ug/L	12/01/25 17:18	VN120125
124-48-1	Dibromochloromethane	0.18	U	1	0.18	1.00	ug/L	12/01/25 17:18	VN120125
106-93-4	1,2-Dibromoethane	0.15	U	1	0.15	1.00	ug/L	12/01/25 17:18	VN120125
127-18-4	Tetrachloroethene	0.23	U	1	0.23	1.00	ug/L	12/01/25 17:18	VN120125
108-90-7	Chlorobenzene	0.12	U	1	0.12	1.00	ug/L	12/01/25 17:18	VN120125
100-41-4	Ethyl Benzene	0.13	U	1	0.13	1.00	ug/L	12/01/25 17:18	VN120125
179601-23-1	m/p-Xylenes	0.24	U	1	0.24	2.00	ug/L	12/01/25 17:18	VN120125
95-47-6	o-Xylene	0.12	U	1	0.12	1.00	ug/L	12/01/25 17:18	VN120125
100-42-5	Styrene	0.15	U	1	0.15	1.00	ug/L	12/01/25 17:18	VN120125
75-25-2	Bromoform	0.19	U	1	0.19	1.00	ug/L	12/01/25 17:18	VN120125



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

Report of Analysis

Client: Remington & Vernick
Project: Saddler Property
Client Sample ID: TB
Lab Sample ID: Q3743-03
Analytical Method: 8260D
Sample Wt/Vol: 5 mL

Level: LOW
Final Vol: 5000 uL

Date Collected: 11/26/25
Date Received: 11/26/25
SDG No.: Q3743
Matrix: Water
% Solid: 0
Test: VOC-TCLVOA-10

CAS Number	Parameter	Conc.	Qua.	DF	MDL	LOQ / CRQL	Units	Date Ana.	BatchID
98-82-8	Isopropylbenzene	0.12	U	1	0.12	1.00	ug/L	12/01/25 17:18	VN120125
79-34-5	1,1,2,2-Tetrachloroethane	0.26	U	1	0.26	1.00	ug/L	12/01/25 17:18	VN120125
541-73-1	1,3-Dichlorobenzene	0.16	U	1	0.16	1.00	ug/L	12/01/25 17:18	VN120125
106-46-7	1,4-Dichlorobenzene	0.19	U	1	0.19	1.00	ug/L	12/01/25 17:18	VN120125
95-50-1	1,2-Dichlorobenzene	0.16	U	1	0.16	1.00	ug/L	12/01/25 17:18	VN120125
96-12-8	1,2-Dibromo-3-Chloropropane	0.53	U	1	0.53	1.00	ug/L	12/01/25 17:18	VN120125
120-82-1	1,2,4-Trichlorobenzene	0.20	U	1	0.20	1.00	ug/L	12/01/25 17:18	VN120125
87-61-6	1,2,3-Trichlorobenzene	0.20	U	1	0.20	1.00	ug/L	12/01/25 17:18	VN120125

SURROGATES

17060-07-0	1,2-Dichloroethane-d4	60.1			74 - 125	120%	SPK: 50
1868-53-7	Dibromofluoromethane	52.4			75 - 124	105%	SPK: 50
2037-26-5	Toluene-d8	50.2			86 - 113	100%	SPK: 50
460-00-4	4-Bromofluorobenzene	53.5			77 - 121	107%	SPK: 50

INTERNAL STANDARDS

INTERNAL STANDARDS		Area Count
363-72-4	Pentafluorobenzene	178000
540-36-3	1,4-Difluorobenzene	408000
3114-55-4	Chlorobenzene-d5	395000
3855-82-1	1,4-Dichlorobenzene-d4	183000

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\VN120125\
 Data File : VN088389.D
 Acq On : 01 Dec 2025 17:18
 Operator : JC\MD
 Sample : Q3743-03
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 TB

Quant Time: Dec 02 00:17:47 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	177560	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.083	114	408212	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.841	117	395290	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.771	152	183327	50.000	ug/l	0.00

System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.559	65	200978	60.085	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	120.180%
35) Dibromofluoromethane	8.147	113	148235	52.400	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	104.800%
50) Toluene-d8	10.541	98	527906	50.154	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	100.300%
62) 4-Bromofluorobenzene	12.829	95	193240	53.506	ug/l	0.00
Spiked Amount	50.000	Range	77 - 121	Recovery	=	107.020%

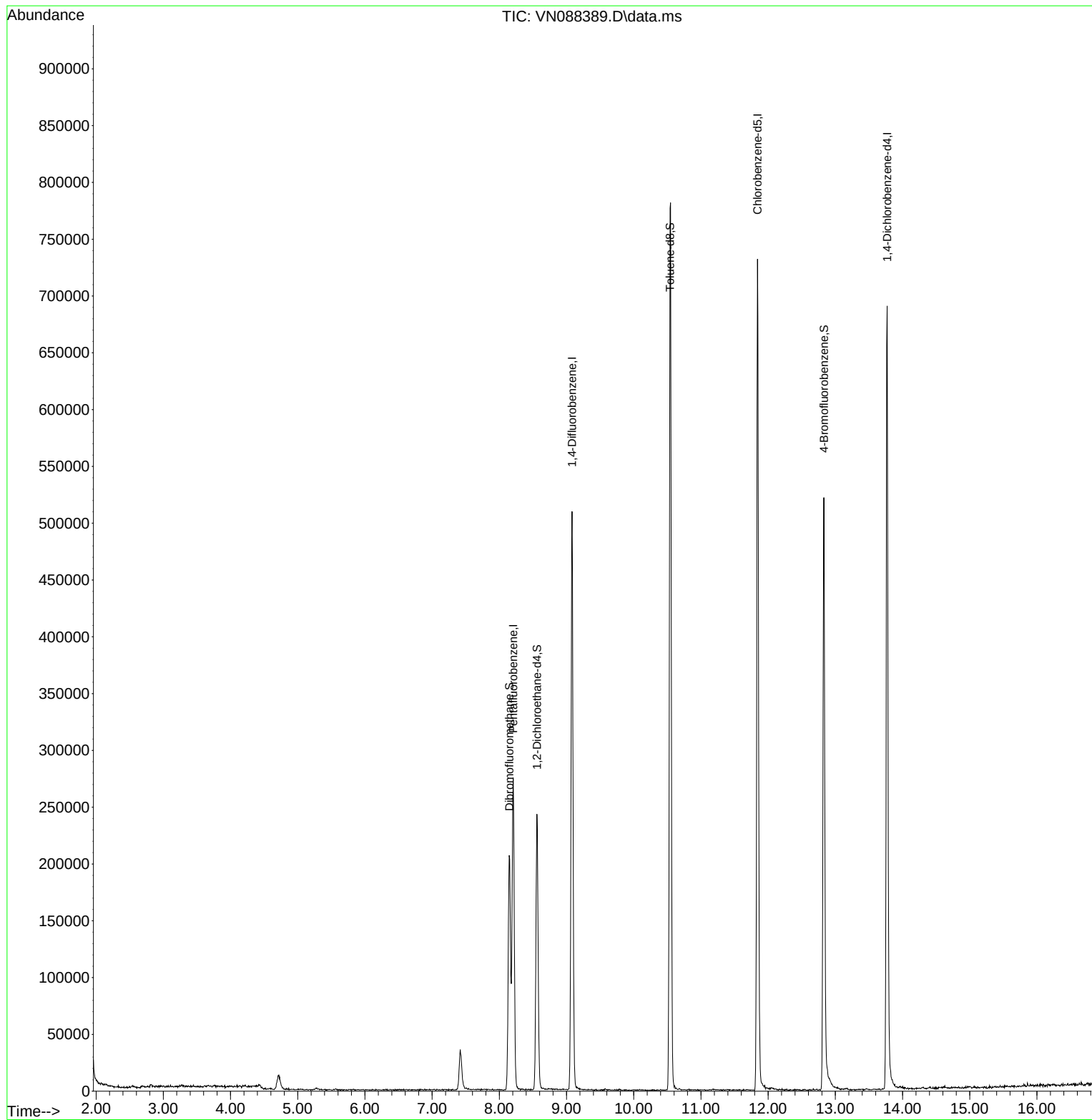
Target Compounds	Qvalue

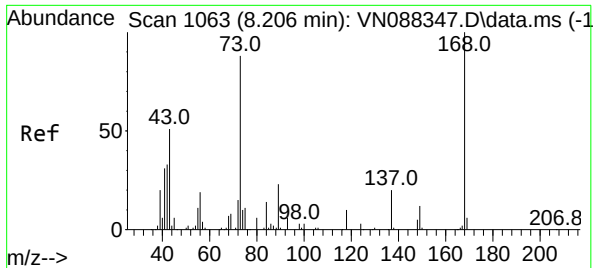
(#) = qualifier out of range (m) = manual integration (+) = signals summed

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN120125\
Data File : VN088389.D
Acq On : 01 Dec 2025 17:18
Operator : JC\MD
Sample : Q3743-03
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
TB

Quant Time: Dec 02 00:17:47 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

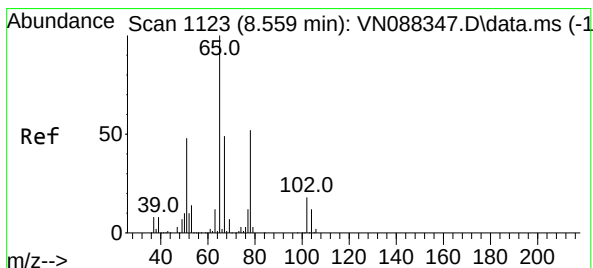
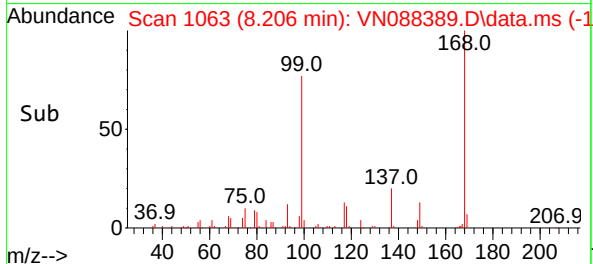
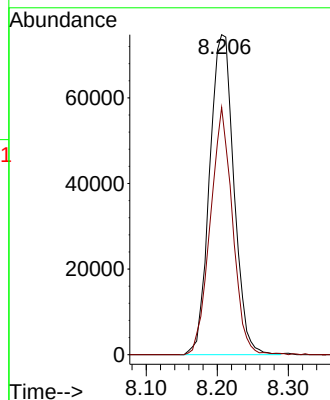
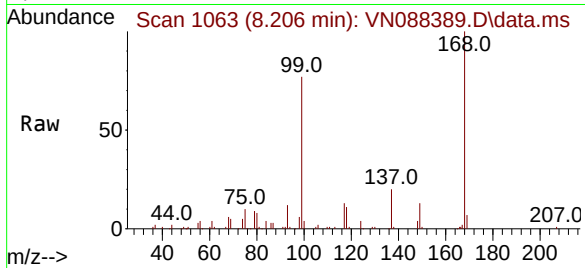




#1
Pentafluorobenzene
Concen: 50.000 ug/l
RT: 8.206 min Scan# 1063
Delta R.T. 0.000 min
Lab File: VN088389.D
Acq: 01 Dec 2025 17:18

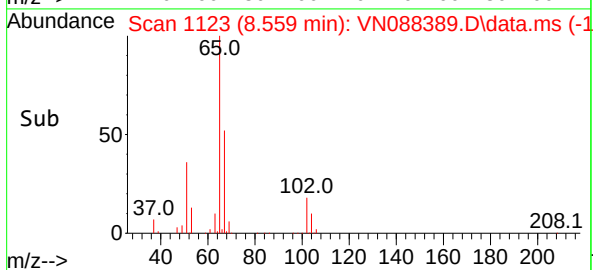
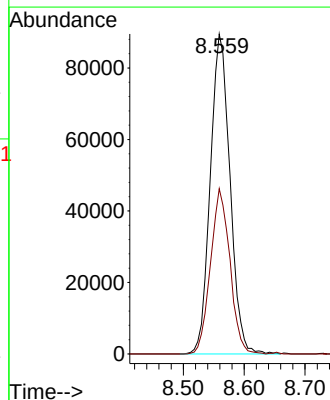
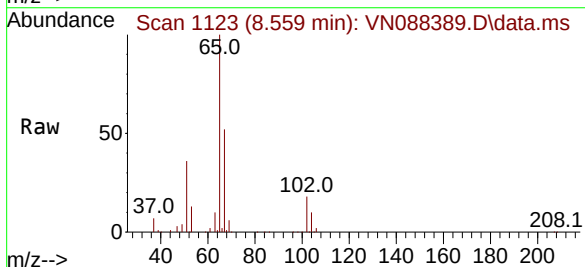
Instrument :
MSVOA_N
ClientSampleId :
TB

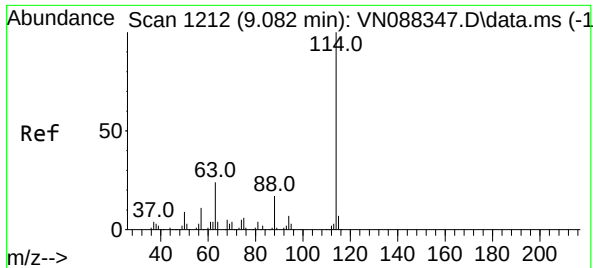
Tgt Ion:168 Resp: 177560
Ion Ratio Lower Upper
168 100
99 77.3 57.1 85.7



#33
1,2-Dichloroethane-d4
Concen: 60.085 ug/l
RT: 8.559 min Scan# 1123
Delta R.T. 0.000 min
Lab File: VN088389.D
Acq: 01 Dec 2025 17:18

Tgt Ion: 65 Resp: 200978
Ion Ratio Lower Upper
65 100
67 50.9 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: VN088389.D

Acq: 01 Dec 2025 17:18

Instrument :

MSVOA_N

ClientSampleId :

TB

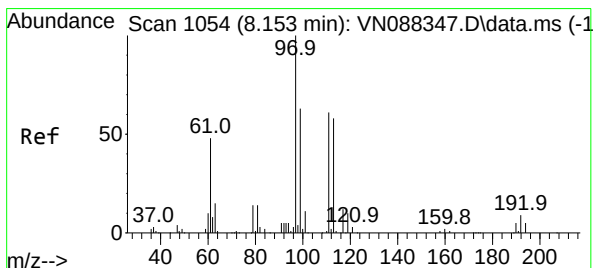
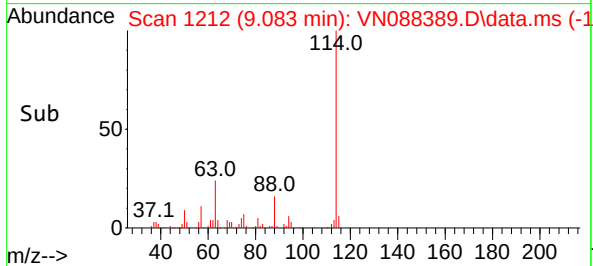
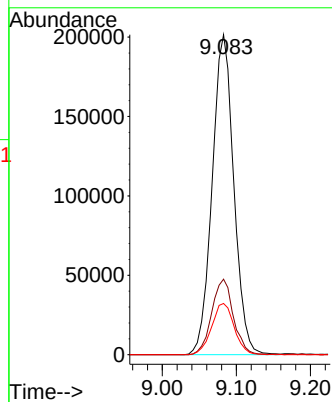
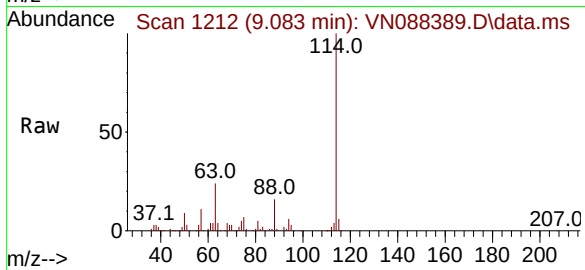
Tgt Ion:114 Resp: 408212

Ion Ratio Lower Upper

114 100

63 23.6 0.0 49.0

88 16.0 0.0 34.0



#35

Dibromofluoromethane

Concen: 52.400 ug/l

RT: 8.147 min Scan# 1053

Delta R.T. -0.006 min

Lab File: VN088389.D

Acq: 01 Dec 2025 17:18

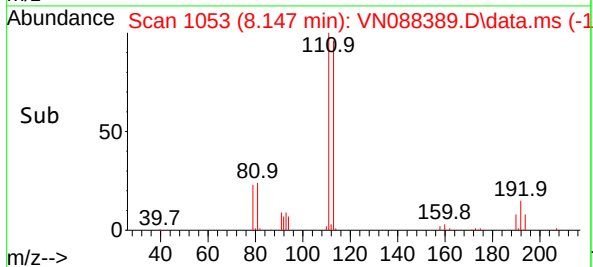
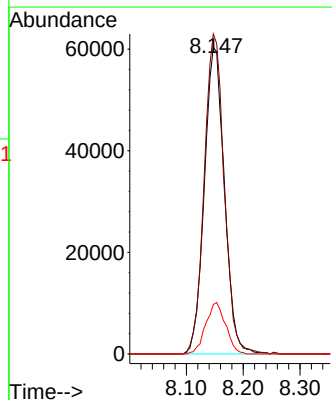
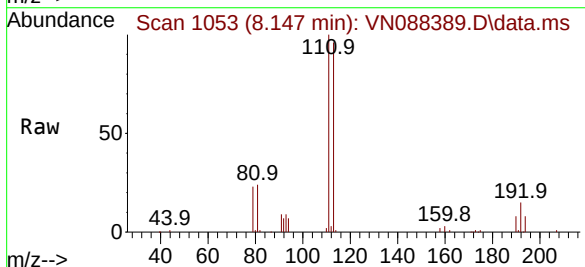
Tgt Ion:113 Resp: 148235

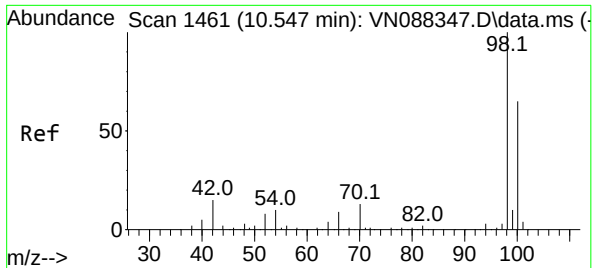
Ion Ratio Lower Upper

113 100

111 103.3 82.5 123.7

192 16.6 12.8 19.2

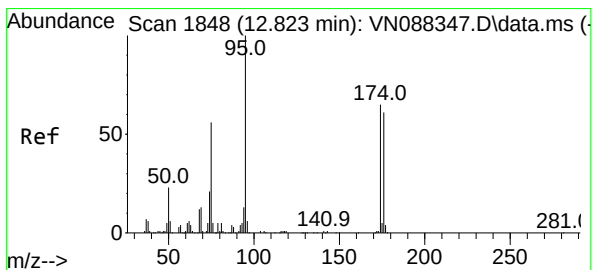
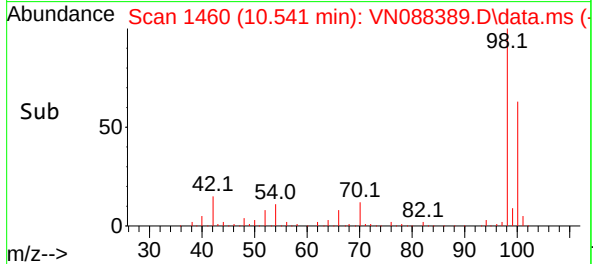
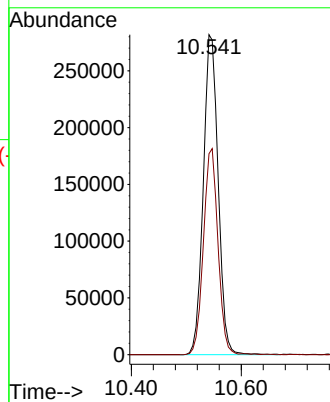
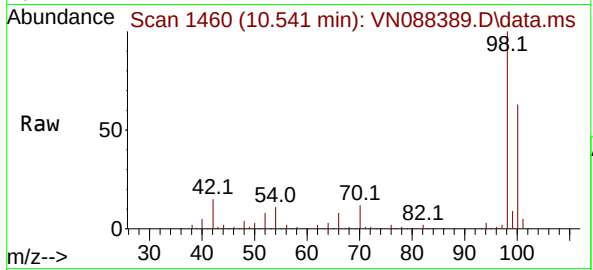




#50
Toluene-d8
Concen: 50.154 ug/l
RT: 10.541 min Scan# 1460
Delta R.T. -0.006 min
Lab File: VN088389.D
Acq: 01 Dec 2025 17:18

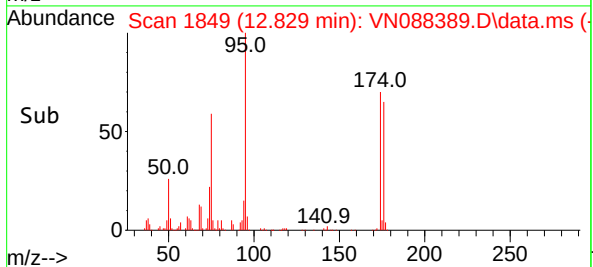
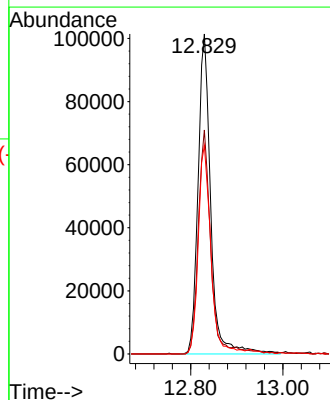
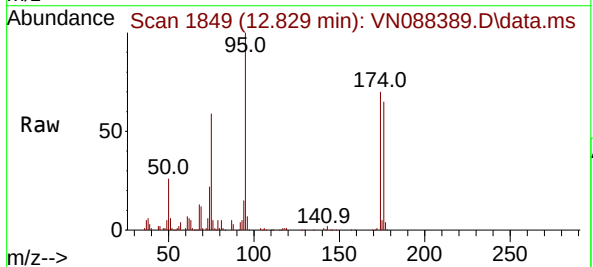
Instrument :
MSVOA_N
ClientSampleId :
TB

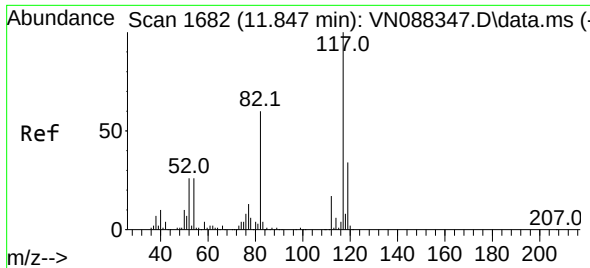
Tgt Ion: 98 Resp: 527906
Ion Ratio Lower Upper
98 100
100 63.3 52.2 78.2



#62
4-Bromofluorobenzene
Concen: 53.506 ug/l
RT: 12.829 min Scan# 1849
Delta R.T. 0.006 min
Lab File: VN088389.D
Acq: 01 Dec 2025 17:18

Tgt Ion: 95 Resp: 193240
Ion Ratio Lower Upper
95 100
174 69.6 0.0 139.0
176 64.8 0.0 132.4

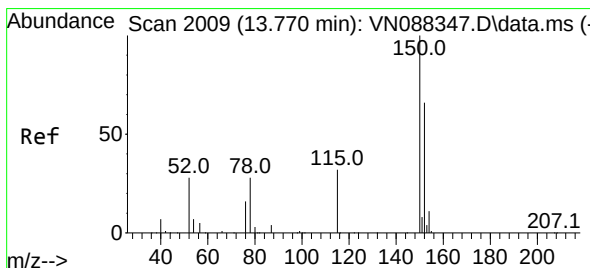
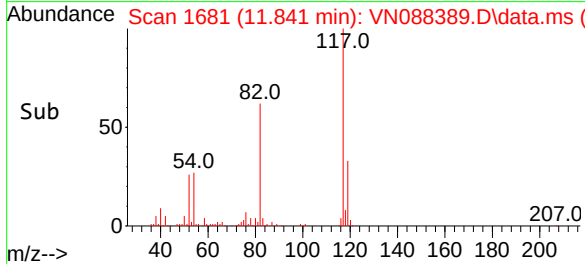
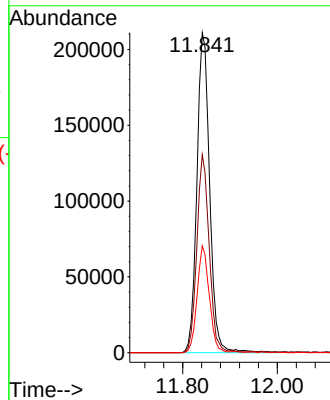
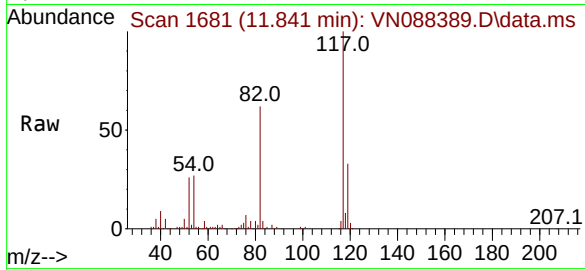




#63
Chlorobenzene-d5
Concen: 50.000 ug/l
RT: 11.841 min Scan# 117
Delta R.T. -0.006 min
Lab File: VN088389.D
Acq: 01 Dec 2025 17:18

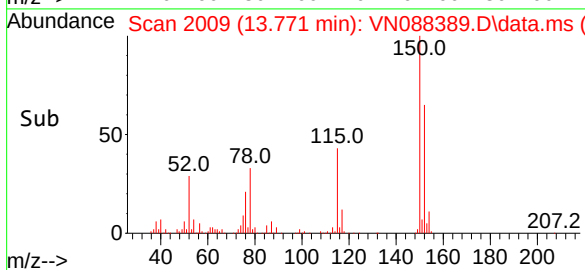
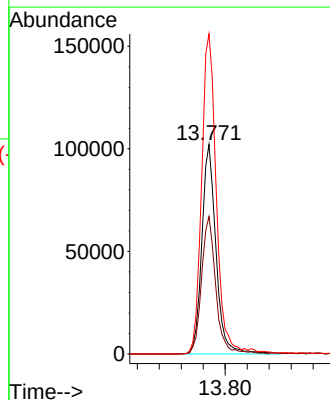
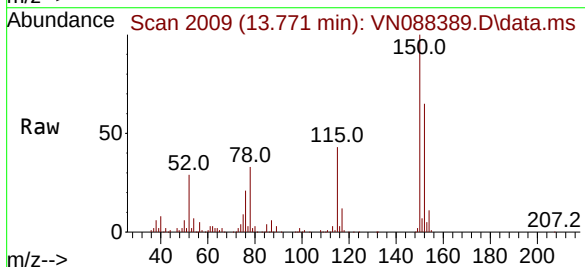
Instrument :
MSVOA_N
ClientSampleId :
TB

Tgt Ion	Ratio	Lower	Upper
117	100		
82	61.9	47.8	71.8
119	33.4	26.9	40.3



#72
1,4-Dichlorobenzene-d4
Concen: 50.000 ug/l
RT: 13.771 min Scan# 2009
Delta R.T. 0.000 min
Lab File: VN088389.D
Acq: 01 Dec 2025 17:18

Tgt Ion	Ratio	Lower	Upper
152	100		
115	65.9	33.0	98.9
150	159.1	0.0	349.0



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN120125\
Data File : VN088389.D
Acq On : 01 Dec 2025 17:18
Operator : JC\MD
Sample : Q3743-03
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
TB

Integration Parameters: RTEINT.P

Integrator: RTE

Smoothing : ON

Sampling : 1

Start Thrs: 0.2

Stop Thrs : 0

Filtering: 5

Min Area: 3 % of largest Peak

Max Peaks: 100

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

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

Title : SW846 8260

Signal : TIC: VN088389.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.718	458	470	479	rBV	12890	45484	3.09%	0.571%
2	7.418	920	929	941	rBV2	35362	97376	6.61%	1.222%
3	8.147	1043	1053	1058	rBV2	206507	512012	34.77%	6.426%
4	8.206	1058	1063	1075	rVB	274076	612264	41.57%	7.684%
5	8.559	1112	1123	1136	rBV	242657	552351	37.51%	6.932%
6	9.083	1203	1212	1224	rBV	508503	1054940	71.63%	13.240%
7	10.547	1452	1461	1470	rBV	781085	1472728	100.00%	18.483%
8	11.841	1673	1681	1698	rBV	731892	1349431	91.63%	16.936%
9	12.829	1840	1849	1868	rBV	521868	998953	67.83%	12.537%
10	13.771	2001	2009	2027	rBV	689035	1272385	86.40%	15.969%

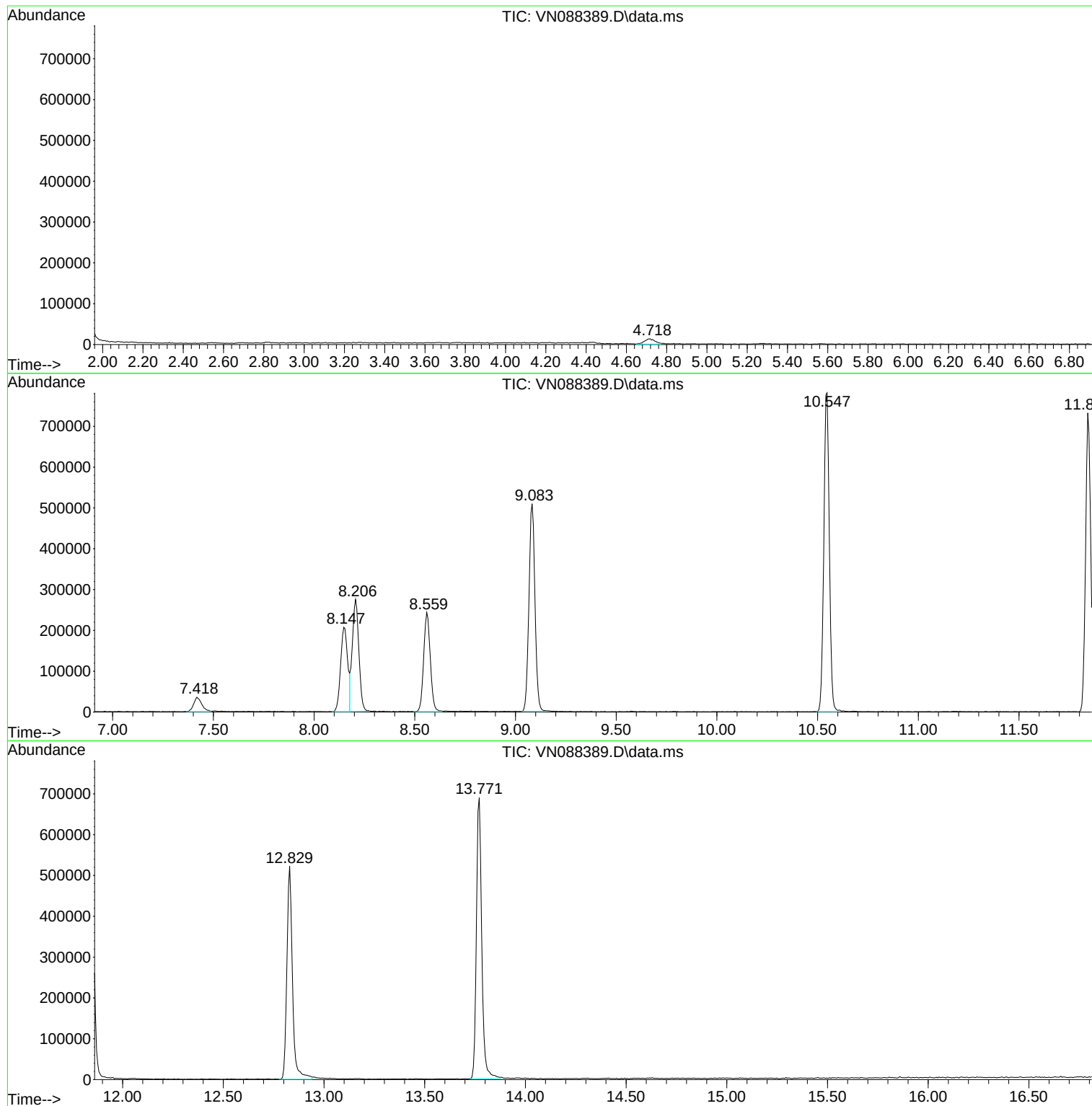
Sum of corrected areas: 7967924

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN120125\
Data File : VN088389.D
Acq On : 01 Dec 2025 17:18
Operator : JC\MD
Sample : Q3743-03
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
TB

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N112425W.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN120125\
Data File : VN088389.D
Acq On : 01 Dec 2025 17:18
Operator : JC\MD
Sample : Q3743-03
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
TB

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N112425W.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

No Library Search Compounds Detected

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN120125\
Data File : VN088389.D
Acq On : 01 Dec 2025 17:18
Operator : JC\MD
Sample : Q3743-03
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
TB

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N112425W.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc

Report of Analysis

Client: Remington & Vernick
Project: Saddler Property
Client Sample ID: TB
Lab Sample ID: Q3743-04
Analytical Method: 8260D
Sample Wt/Vol: 5 mL

Level: LOW
Final Vol: 5000 uL

Date Collected: 11/26/25
Date Received: 11/26/25
SDG No.: Q3743
Matrix: Water
% Solid: 0
Test: VOC-TCLVOA-10

CAS Number	Parameter	Conc.	Qua.	DF	MDL	LOQ / CRQL	Units	Date Ana.	BatchID
TARGETS									
75-71-8	Dichlorodifluoromethane	0.22	U	1	0.22	1.00	ug/L	12/01/25 17:39	VN120125
74-87-3	Chloromethane	0.32	U	1	0.32	1.00	ug/L	12/01/25 17:39	VN120125
75-01-4	Vinyl Chloride	0.26	U	1	0.26	1.00	ug/L	12/01/25 17:39	VN120125
74-83-9	Bromomethane	1.40	U	1	1.40	5.00	ug/L	12/01/25 17:39	VN120125
75-00-3	Chloroethane	0.47	U	1	0.47	1.00	ug/L	12/01/25 17:39	VN120125
75-69-4	Trichlorofluoromethane	0.33	U	1	0.33	1.00	ug/L	12/01/25 17:39	VN120125
76-13-1	1,1,2-Trichlorotrifluoroethane	0.25	U	1	0.25	1.00	ug/L	12/01/25 17:39	VN120125
75-35-4	1,1-Dichloroethene	0.23	U	1	0.23	1.00	ug/L	12/01/25 17:39	VN120125
67-64-1	Acetone	1.50	U	1	1.50	5.00	ug/L	12/01/25 17:39	VN120125
75-15-0	Carbon Disulfide	0.21	U	1	0.21	1.00	ug/L	12/01/25 17:39	VN120125
1634-04-4	Methyl tert-butyl Ether	0.16	U	1	0.16	1.00	ug/L	12/01/25 17:39	VN120125
79-20-9	Methyl Acetate	0.27	U	1	0.27	1.00	ug/L	12/01/25 17:39	VN120125
75-09-2	Methylene Chloride	0.28	U	1	0.28	1.00	ug/L	12/01/25 17:39	VN120125
156-60-5	trans-1,2-Dichloroethene	0.23	U	1	0.23	1.00	ug/L	12/01/25 17:39	VN120125
75-34-3	1,1-Dichloroethane	0.23	U	1	0.23	1.00	ug/L	12/01/25 17:39	VN120125
110-82-7	Cyclohexane	1.50	U	1	1.50	5.00	ug/L	12/01/25 17:39	VN120125
78-93-3	2-Butanone	0.98	U	1	0.98	5.00	ug/L	12/01/25 17:39	VN120125
56-23-5	Carbon Tetrachloride	0.25	U	1	0.25	1.00	ug/L	12/01/25 17:39	VN120125
156-59-2	cis-1,2-Dichloroethene	0.19	U	1	0.19	1.00	ug/L	12/01/25 17:39	VN120125
74-97-5	Bromochloromethane	0.22	U	1	0.22	1.00	ug/L	12/01/25 17:39	VN120125
67-66-3	Chloroform	0.25	U	1	0.25	1.00	ug/L	12/01/25 17:39	VN120125
71-55-6	1,1,1-Trichloroethane	0.20	U	1	0.20	1.00	ug/L	12/01/25 17:39	VN120125
108-87-2	Methylcyclohexane	0.16	U	1	0.16	1.00	ug/L	12/01/25 17:39	VN120125
71-43-2	Benzene	0.15	U	1	0.15	1.00	ug/L	12/01/25 17:39	VN120125
107-06-2	1,2-Dichloroethane	0.22	U	1	0.22	1.00	ug/L	12/01/25 17:39	VN120125
79-01-6	Trichloroethene	0.090	U	1	0.090	1.00	ug/L	12/01/25 17:39	VN120125
78-87-5	1,2-Dichloropropane	0.20	U	1	0.20	1.00	ug/L	12/01/25 17:39	VN120125
75-27-4	Bromodichloromethane	0.22	U	1	0.22	1.00	ug/L	12/01/25 17:39	VN120125
108-10-1	4-Methyl-2-Pentanone	0.68	U	1	0.68	5.00	ug/L	12/01/25 17:39	VN120125
108-88-3	Toluene	0.14	U	1	0.14	1.00	ug/L	12/01/25 17:39	VN120125
10061-02-6	t-1,3-Dichloropropene	0.17	U	1	0.17	1.00	ug/L	12/01/25 17:39	VN120125
10061-01-5	cis-1,3-Dichloropropene	0.16	U	1	0.16	1.00	ug/L	12/01/25 17:39	VN120125
79-00-5	1,1,2-Trichloroethane	0.21	U	1	0.21	1.00	ug/L	12/01/25 17:39	VN120125
591-78-6	2-Hexanone	0.89	U	1	0.89	5.00	ug/L	12/01/25 17:39	VN120125
124-48-1	Dibromochloromethane	0.18	U	1	0.18	1.00	ug/L	12/01/25 17:39	VN120125
106-93-4	1,2-Dibromoethane	0.15	U	1	0.15	1.00	ug/L	12/01/25 17:39	VN120125
127-18-4	Tetrachloroethene	0.23	U	1	0.23	1.00	ug/L	12/01/25 17:39	VN120125
108-90-7	Chlorobenzene	0.12	U	1	0.12	1.00	ug/L	12/01/25 17:39	VN120125
100-41-4	Ethyl Benzene	0.13	U	1	0.13	1.00	ug/L	12/01/25 17:39	VN120125
179601-23-1	m/p-Xylenes	0.24	U	1	0.24	2.00	ug/L	12/01/25 17:39	VN120125
95-47-6	o-Xylene	0.12	U	1	0.12	1.00	ug/L	12/01/25 17:39	VN120125
100-42-5	Styrene	0.15	U	1	0.15	1.00	ug/L	12/01/25 17:39	VN120125
75-25-2	Bromoform	0.19	U	1	0.19	1.00	ug/L	12/01/25 17:39	VN120125



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

Report of Analysis

Client: Remington & Vernick
Project: Saddler Property
Client Sample ID: TB
Lab Sample ID: Q3743-04
Analytical Method: 8260D
Sample Wt/Vol: 5 mL

Level: LOW
Final Vol: 5000 uL

Date Collected: 11/26/25
Date Received: 11/26/25
SDG No.: Q3743
Matrix: Water
% Solid: 0
Test: VOC-TCLVOA-10

CAS Number	Parameter	Conc.	Qua.	DF	MDL	LOQ / CRQL	Units	Date Ana.	BatchID
98-82-8	Isopropylbenzene	0.12	U	1	0.12	1.00	ug/L	12/01/25 17:39	VN120125
79-34-5	1,1,2,2-Tetrachloroethane	0.26	U	1	0.26	1.00	ug/L	12/01/25 17:39	VN120125
541-73-1	1,3-Dichlorobenzene	0.16	U	1	0.16	1.00	ug/L	12/01/25 17:39	VN120125
106-46-7	1,4-Dichlorobenzene	0.19	U	1	0.19	1.00	ug/L	12/01/25 17:39	VN120125
95-50-1	1,2-Dichlorobenzene	0.16	U	1	0.16	1.00	ug/L	12/01/25 17:39	VN120125
96-12-8	1,2-Dibromo-3-Chloropropane	0.53	U	1	0.53	1.00	ug/L	12/01/25 17:39	VN120125
120-82-1	1,2,4-Trichlorobenzene	0.20	U	1	0.20	1.00	ug/L	12/01/25 17:39	VN120125
87-61-6	1,2,3-Trichlorobenzene	0.20	U	1	0.20	1.00	ug/L	12/01/25 17:39	VN120125

SURROGATES

17060-07-0	1,2-Dichloroethane-d4	60.9			74 - 125	122%	SPK: 50
1868-53-7	Dibromofluoromethane	53.0			75 - 124	106%	SPK: 50
2037-26-5	Toluene-d8	49.8			86 - 113	100%	SPK: 50
460-00-4	4-Bromofluorobenzene	53.2			77 - 121	106%	SPK: 50

INTERNAL STANDARDS

INTERNAL STANDARDS		Area Count
363-72-4	Pentafluorobenzene	168000
540-36-3	1,4-Difluorobenzene	387000
3114-55-4	Chlorobenzene-d5	373000
3855-82-1	1,4-Dichlorobenzene-d4	176000

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\VN120125\
 Data File : VN088390.D
 Acq On : 01 Dec 2025 17:39
 Operator : JC\MD
 Sample : Q3743-04
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 TB

Quant Time: Dec 02 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	168109	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.083	114	386817	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.841	117	372510	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.770	152	175948	50.000	ug/l	0.00

System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.559	65	192789	60.877	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	121.760%
35) Dibromofluoromethane	8.147	113	142167	53.035	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	106.080%
50) Toluene-d8	10.547	98	497084	49.838	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	99.680%
62) 4-Bromofluorobenzene	12.829	95	182162	53.228	ug/l	0.00
Spiked Amount	50.000	Range	77 - 121	Recovery	=	106.460%

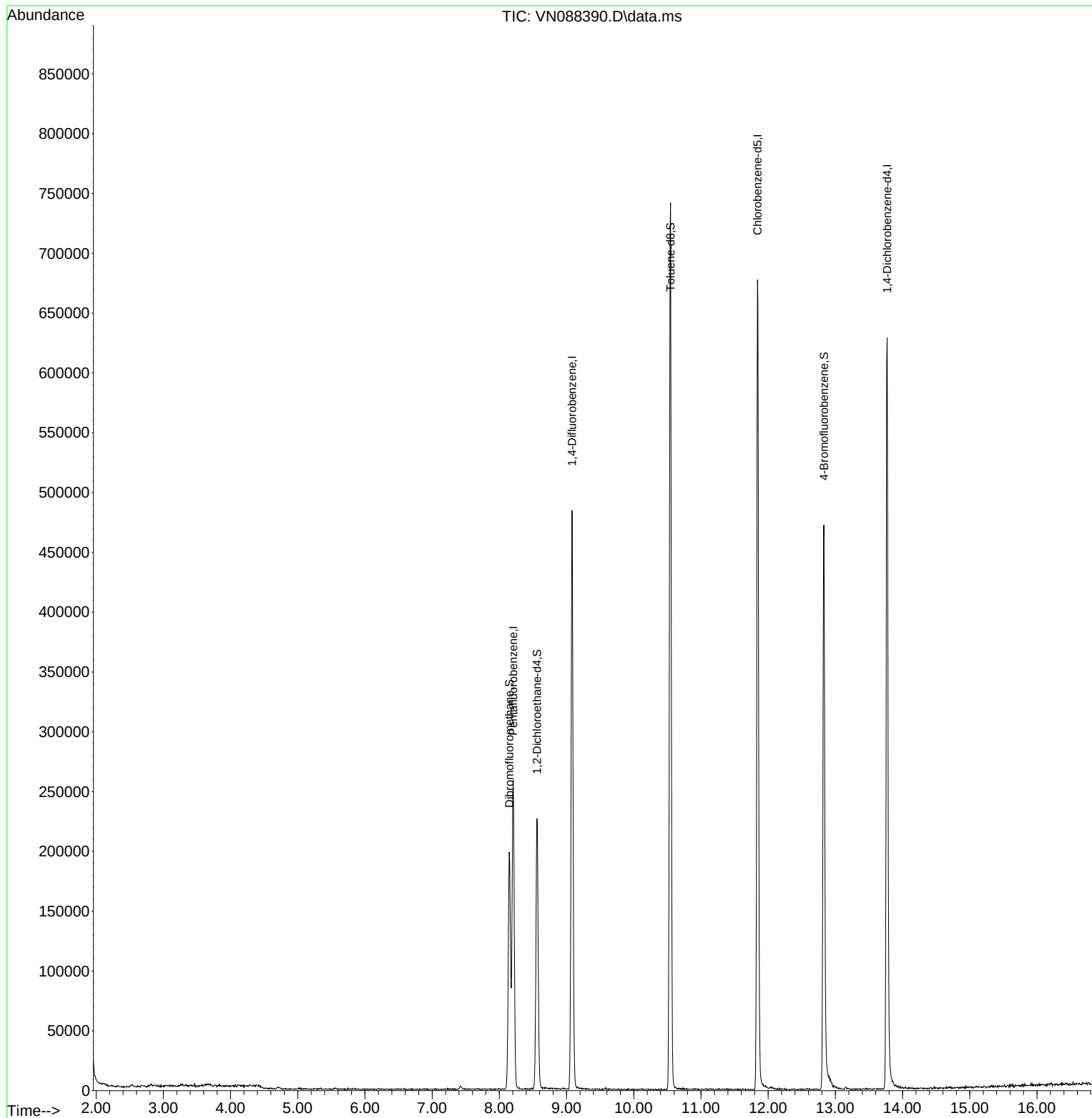
Target Compounds	Qvalue

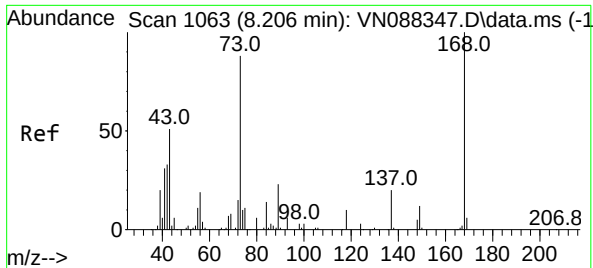
(#) = qualifier out of range (m) = manual integration (+) = signals summed

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN120125\
Data File : VN088390.D
Acq On : 01 Dec 2025 17:39
Operator : JC\MD
Sample : Q3743-04
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
TB

Quant Time: Dec 02 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

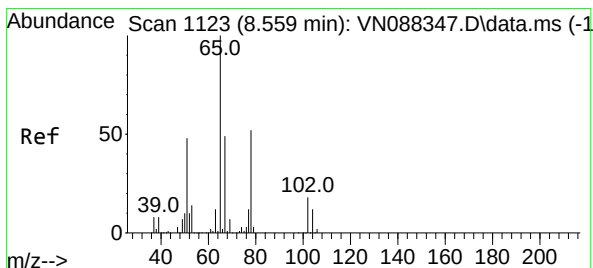
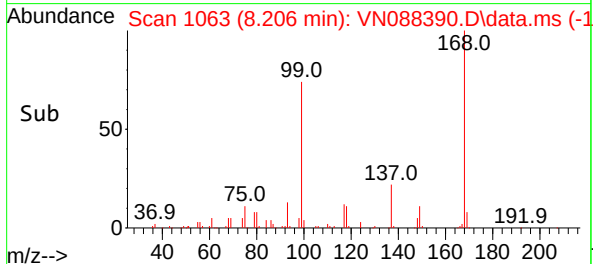
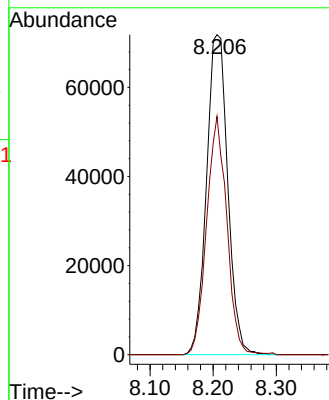
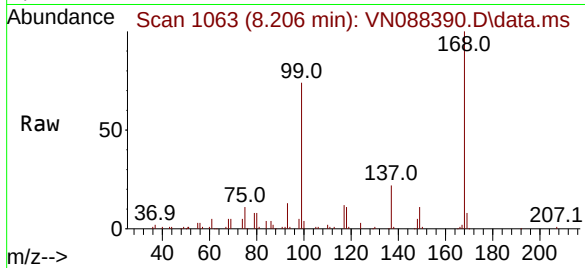




#1
Pentafluorobenzene
Concen: 50.000 ug/l
RT: 8.206 min Scan# 1063
Delta R.T. 0.000 min
Lab File: VN088390.D
Acq: 01 Dec 2025 17:39

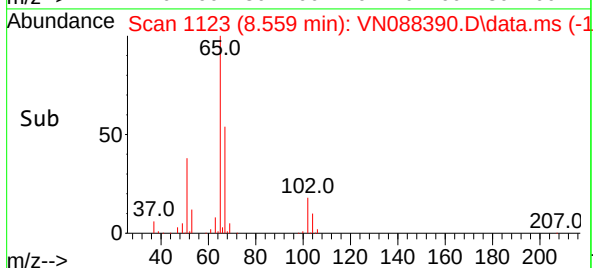
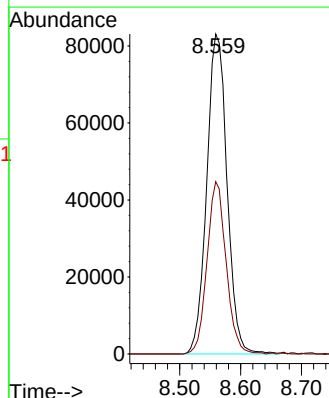
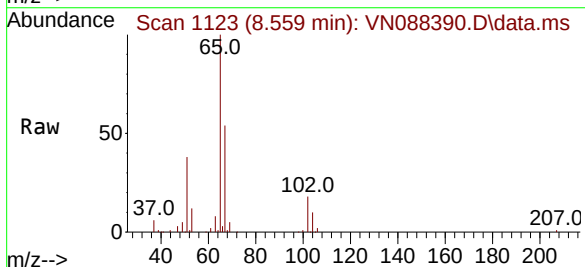
Instrument :
MSVOA_N
ClientSampleId :
TB

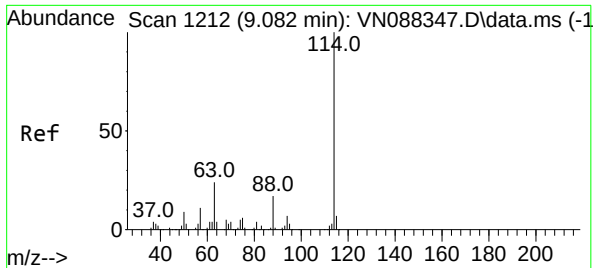
Tgt Ion:168 Resp: 168109
Ion Ratio Lower Upper
168 100
99 74.5 57.1 85.7



#33
1,2-Dichloroethane-d4
Concen: 60.877 ug/l
RT: 8.559 min Scan# 1123
Delta R.T. 0.000 min
Lab File: VN088390.D
Acq: 01 Dec 2025 17:39

Tgt Ion: 65 Resp: 192789
Ion Ratio Lower Upper
65 100
67 50.7 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: VN088390.D

Acq: 01 Dec 2025 17:39

Instrument :

MSVOA_N

ClientSampleId :

TB

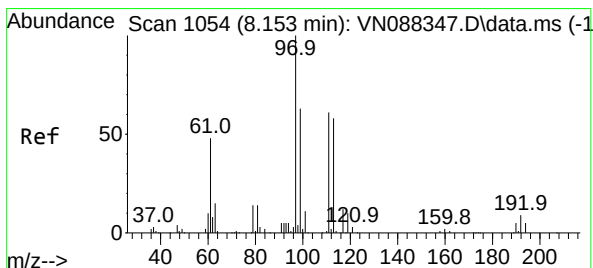
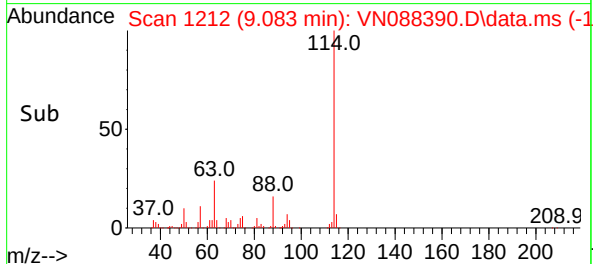
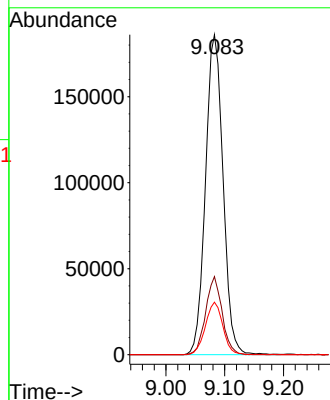
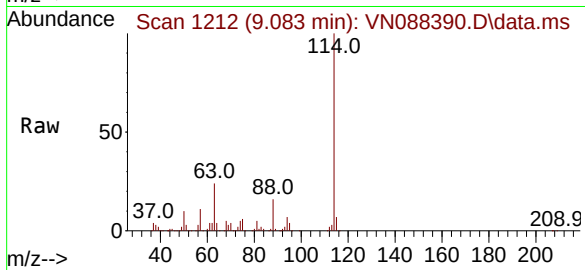
Tgt Ion:114 Resp: 386817

Ion Ratio Lower Upper

114 100

63 24.4 0.0 49.0

88 16.4 0.0 34.0



#35

Dibromofluoromethane

Concen: 53.035 ug/l

RT: 8.147 min Scan# 1053

Delta R.T. -0.006 min

Lab File: VN088390.D

Acq: 01 Dec 2025 17:39

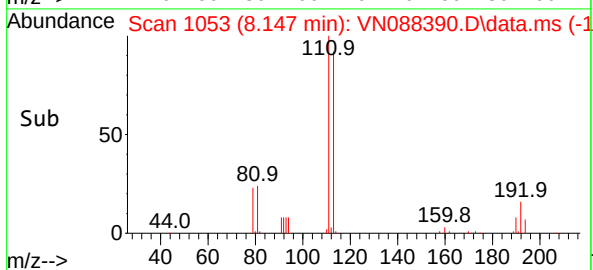
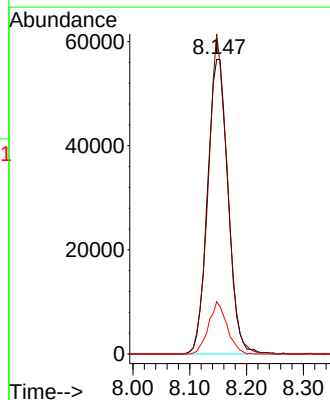
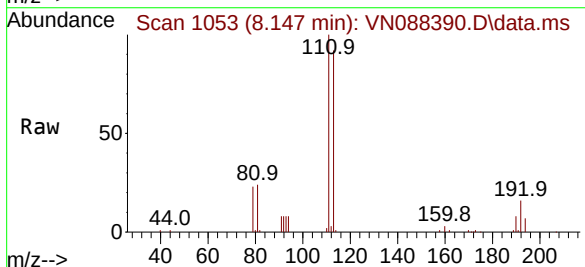
Tgt Ion:113 Resp: 142167

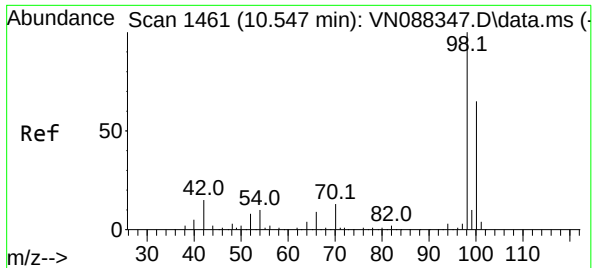
Ion Ratio Lower Upper

113 100

111 100.6 82.5 123.7

192 16.0 12.8 19.2

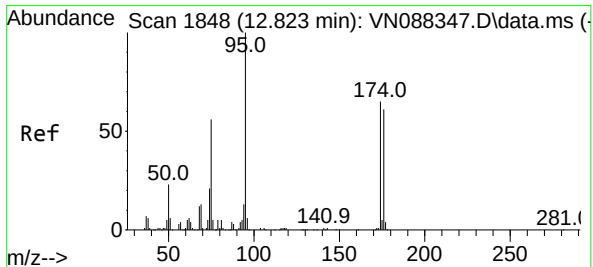
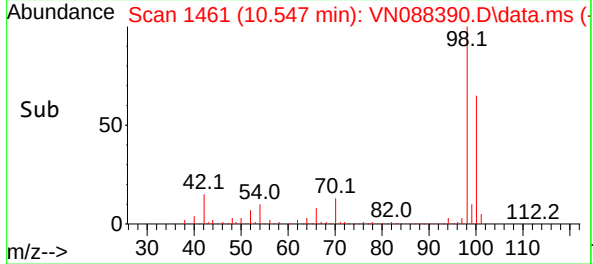
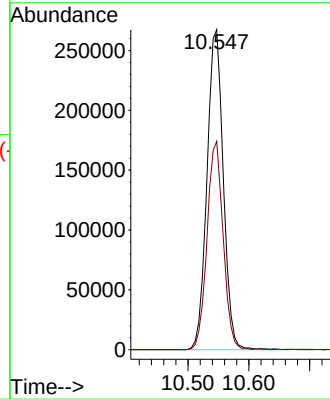
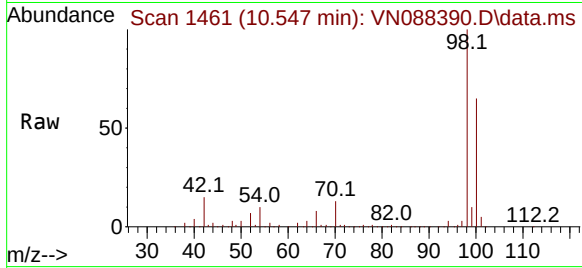




#50
Toluene-d8
Concen: 49.838 ug/l
RT: 10.547 min Scan# 1461
Delta R.T. 0.000 min
Lab File: VN088390.D
Acq: 01 Dec 2025 17:39

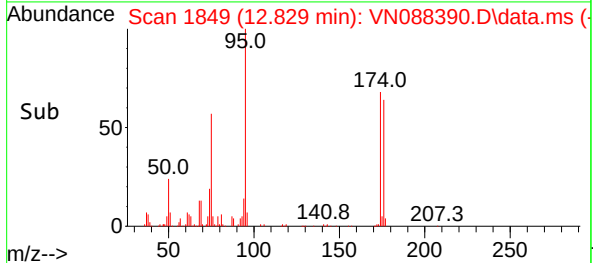
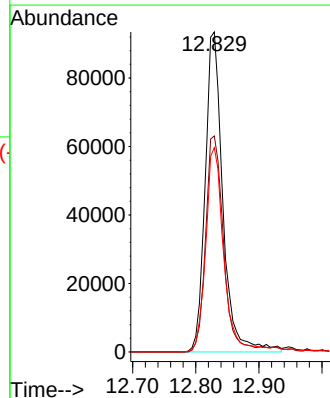
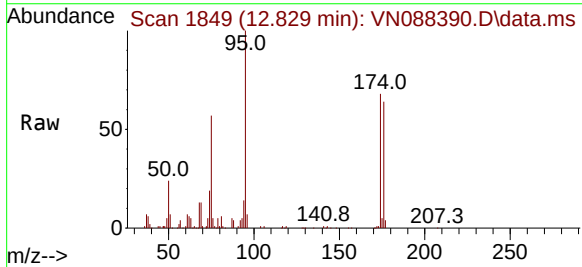
Instrument :
MSVOA_N
ClientSampleId :
TB

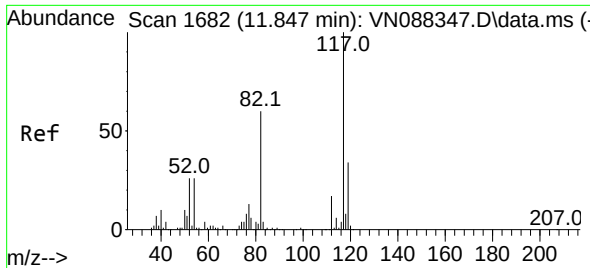
Tgt Ion: 98 Resp: 497084
Ion Ratio Lower Upper
98 100
100 64.3 52.2 78.2



#62
4-Bromofluorobenzene
Concen: 53.228 ug/l
RT: 12.829 min Scan# 1849
Delta R.T. 0.006 min
Lab File: VN088390.D
Acq: 01 Dec 2025 17:39

Tgt Ion: 95 Resp: 182162
Ion Ratio Lower Upper
95 100
174 67.6 0.0 139.0
176 64.8 0.0 132.4

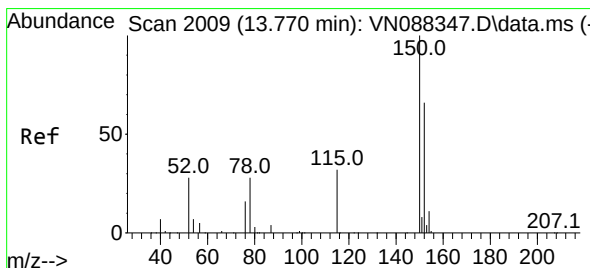
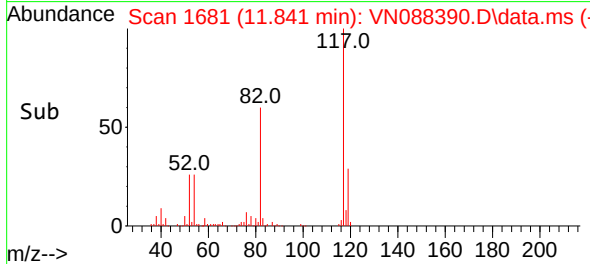
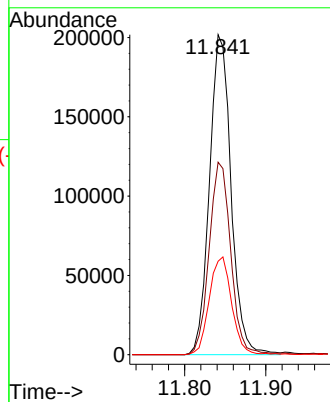
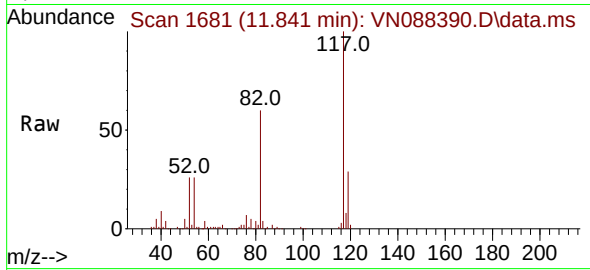




#63
Chlorobenzene-d5
Concen: 50.000 ug/l
RT: 11.841 min Scan# 11
Delta R.T. -0.006 min
Lab File: VN088390.D
Acq: 01 Dec 2025 17:39

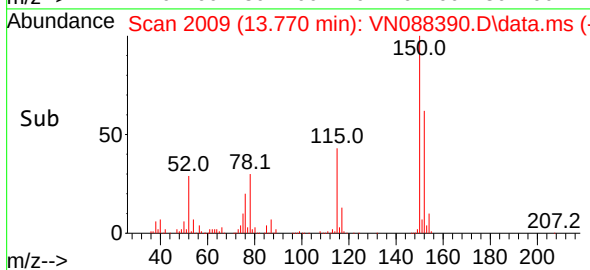
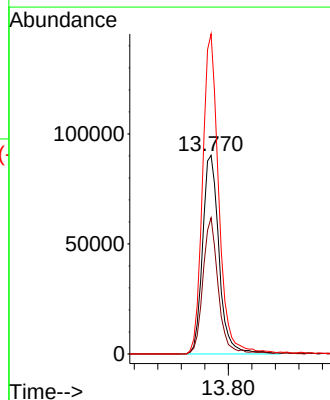
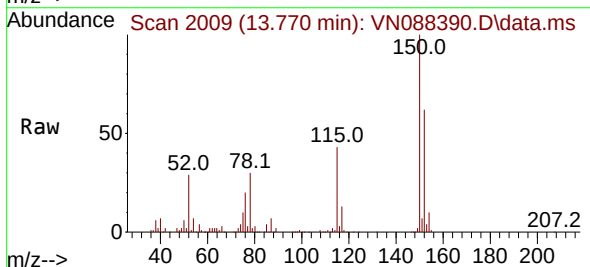
Instrument :
MSVOA_N
ClientSampleId :
TB

Tgt Ion	Ratio	Lower	Upper
117	100		
82	60.1	47.8	71.8
119	29.2	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: VN088390.D
Acq: 01 Dec 2025 17:39

Tgt Ion	Ratio	Lower	Upper
152	100		
115	63.9	33.0	98.9
150	158.2	0.0	349.0



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN120125\
Data File : VN088390.D
Acq On : 01 Dec 2025 17:39
Operator : JC\MD
Sample : Q3743-04
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
TB

Integration Parameters: RTEINT.P

Integrator: RTE

Smoothing : ON

Filtering: 5

Sampling : 1

Min Area: 3 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

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

Title : SW846 8260

Signal : TIC: VN088390.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	8.147	1041	1053	1058	rBV	198377	478562	34.38%	6.522%
2	8.206	1058	1063	1075	rVB	257933	574139	41.25%	7.824%
3	8.559	1114	1123	1135	rBV	226331	521824	37.49%	7.111%
4	9.083	1203	1212	1224	rBV	483945	990791	71.18%	13.503%
5	10.547	1452	1461	1472	rBV	741514	1391998	100.00%	18.970%
6	11.841	1673	1681	1694	rBV	677076	1270448	91.27%	17.314%
7	12.829	1836	1849	1861	rBV	472178	913206	65.60%	12.445%
8	13.770	1999	2009	2022	rBV	628243	1196831	85.98%	16.310%

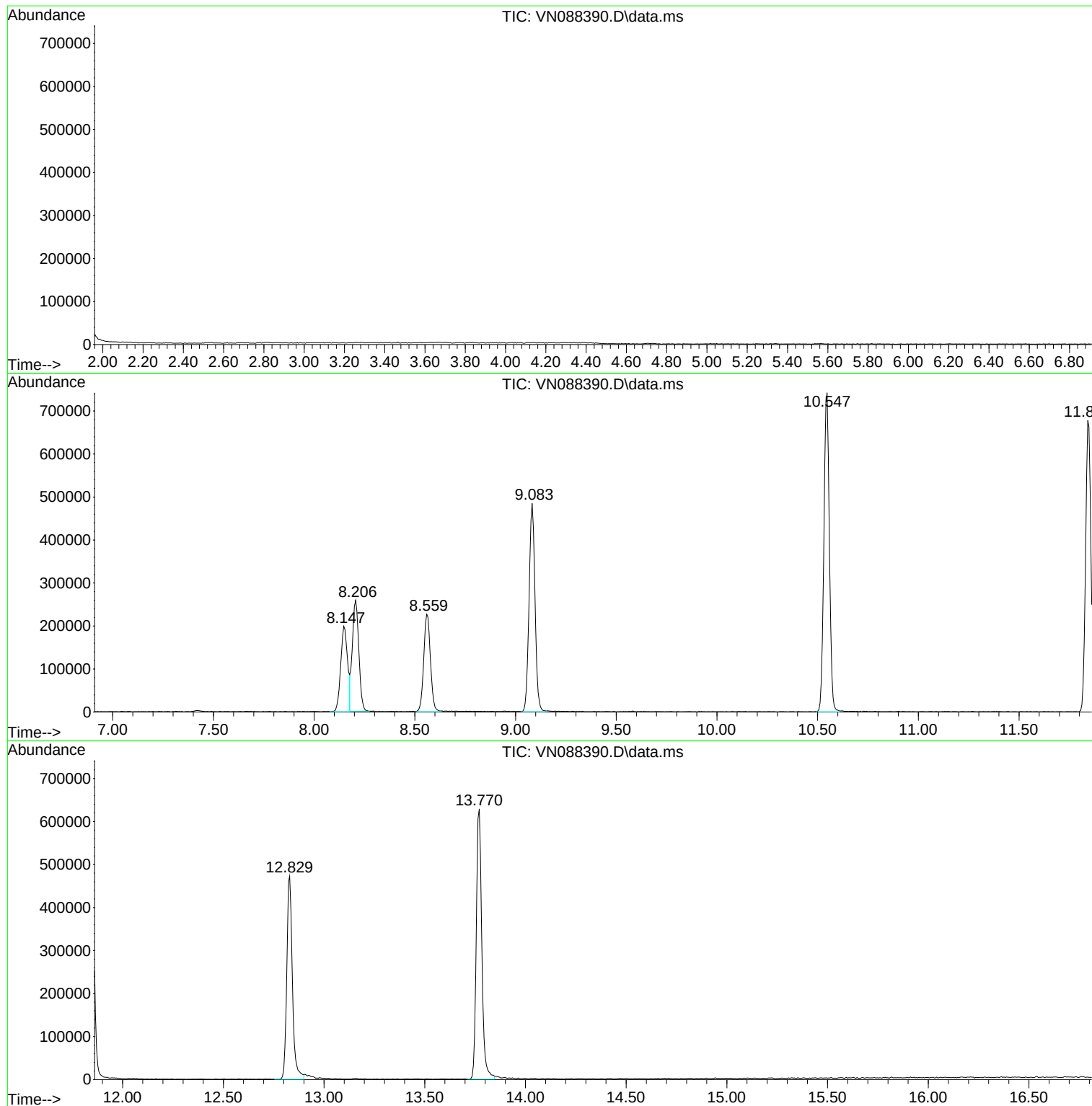
Sum of corrected areas: 7337799

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN120125\
Data File : VN088390.D
Acq On : 01 Dec 2025 17:39
Operator : JC\MD
Sample : Q3743-04
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
TB

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N112425W.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN120125\
Data File : VN088390.D
Acq On : 01 Dec 2025 17:39
Operator : JC\MD
Sample : Q3743-04
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
TB

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N112425W.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

No Library Search Compounds Detected

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN120125\
Data File : VN088390.D
Acq On : 01 Dec 2025 17:39
Operator : JC\MD
Sample : Q3743-04
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
TB

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N112425W.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc



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: REMI01
 Lab Code: ACE SDG No.: Q3743
 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	
Dichlorodifluoromethane	0.719	0.676	0.796	0.741	0.838	0.736	0.751	7.7	
Chloromethane	0.911	0.800	0.805	0.801	0.759	0.764	0.807	6.8	
Vinyl Chloride	0.804	0.844	0.856	0.812	0.837	0.772	0.821	3.8	
Bromomethane		0.407	0.392	0.396	0.387	0.392	0.395	1.9	
Chloroethane	0.555	0.503	0.496	0.505	0.502	0.485	0.508	4.8	
Trichlorofluoromethane	1.234	1.170	1.168	1.092	1.201	1.055	1.153	5.8	
1,1,2-Trichlorotrifluoroethane	0.660	0.635	0.639	0.564	0.633	0.553	0.614	7.2	
1,1-Dichloroethene	0.722	0.620	0.601	0.568	0.595	0.554	0.610	9.8	
Acetone	0.412	0.355	0.322	0.335	0.305	0.314	0.341	11.5	
Carbon Disulfide	2.061	2.043	2.043	1.983	1.996	1.833	1.993	4.2	
Methyl tert-butyl Ether	2.126	2.239	2.245	2.504	2.265	2.333	2.285	5.5	
Methyl Acetate	1.072	1.048	0.957	1.049	0.972	0.985	1.014	4.7	
Methylene Chloride	0.797	0.752	0.704	0.756	0.686	0.678	0.729	6.4	
trans-1,2-Dichloroethene	0.752	0.668	0.676	0.686	0.647	0.626	0.676	6.4	
1,1-Dichloroethane	1.389	1.431	1.348	1.427	1.321	1.286	1.367	4.3	
Cyclohexane		1.450	1.264	1.133	1.237	1.109	1.239	10.9	
2-Butanone	0.509	0.538	0.548	0.596	0.540	0.556	0.548	5.2	
Carbon Tetrachloride	0.565	0.504	0.548	0.499	0.570	0.509	0.532	6	
cis-1,2-Dichloroethene	0.797	0.801	0.794	0.829	0.747	0.744	0.785	4.2	
Bromochloromethane	0.724	0.722	0.684	0.735	0.624	0.597	0.681	8.5	
Chloroform	1.377	1.362	1.351	1.392	1.291	1.285	1.343	3.3	
1,1,1-Trichloroethane	1.247	1.219	1.192	1.158	1.166	1.104	1.181	4.3	
Methylcyclohexane	0.534	0.542	0.606	0.527	0.662	0.561	0.572	9.1	
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	
1,2-Dichloropropane	0.426	0.402	0.402	0.406	0.396	0.381	0.402	3.6	
Bromodichloromethane	0.602	0.586	0.586	0.589	0.583	0.566	0.586	2	
4-Methyl-2-Pentanone	0.540	0.616	0.661	0.673	0.658	0.651	0.633	7.8	
Toluene	0.829	0.885	0.940	0.913	0.960	0.911	0.906	5.1	

* 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.



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

VOLATILE ORGANICS INITIAL CALIBRATION DATA

Lab Name: Alliance Contract: REMI01
 Lab Code: ACE SDG No.: Q3743
 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	
t-1,3-Dichloropropene	0.567	0.553	0.598	0.631	0.631	0.625	0.601	5.7	
cis-1,3-Dichloropropene	0.590	0.618	0.628	0.662	0.654	0.640	0.632	4.2	
1,1,2-Trichloroethane	0.350	0.359	0.364	0.353	0.341	0.336	0.351	3	
2-Hexanone	0.359	0.301	0.406	0.436	0.442	0.450	0.399	14.7	
Dibromochloromethane	0.384	0.386	0.406	0.409	0.410	0.404	0.400	2.9	
1,2-Dibromoethane	0.304	0.367	0.364	0.369	0.367	0.359	0.355	7.2	
Tetrachloroethene	0.426	0.388	0.413	0.377	0.407	0.367	0.396	5.7	
Chlorobenzene	1.114	1.142	1.161	1.106	1.145	1.070	1.123	3	
Ethyl Benzene	1.881	1.860	2.014	1.938	2.133	1.952	1.963	5.1	
m/p-Xylenes	0.658	0.690	0.769	0.733	0.788	0.728	0.728	6.6	
o-Xylene	0.670	0.619	0.722	0.696	0.749	0.708	0.694	6.5	
Styrene	0.884	1.025	1.163	1.193	1.284	1.210	1.127	12.9	
Bromoform	0.224	0.273	0.285	0.285	0.302	0.295	0.277	10.1	
Isopropylbenzene	3.339	3.528	3.944	3.624	4.040	3.706	3.697	7.1	
1,1,2,2-Tetrachloroethane	1.377	1.218	1.231	1.127	1.135	1.082	1.195	8.9	
1,3-Dichlorobenzene	1.730	1.573	1.720	1.606	1.676	1.574	1.647	4.3	
1,4-Dichlorobenzene	1.747	1.825	1.786	1.681	1.702	1.608	1.725	4.5	
1,2-Dichlorobenzene	1.518	1.574	1.583	1.521	1.595	1.505	1.550	2.5	
1,2-Dibromo-3-Chloropropane	0.284	0.284	0.284	0.288	0.300	0.300	0.290	2.8	
1,2,4-Trichlorobenzene	0.998	0.864	0.901	0.889	0.955	0.917	0.921	5.3	
1,2,3-Trichlorobenzene	0.942	0.850	0.888	0.875	0.930	0.891	0.896	3.8	
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 : VSTDIC001
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 2 Sample Multiplier: 1

Instrument :

MSVOA_N

ClientSampleId :

VSTDIC001

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

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

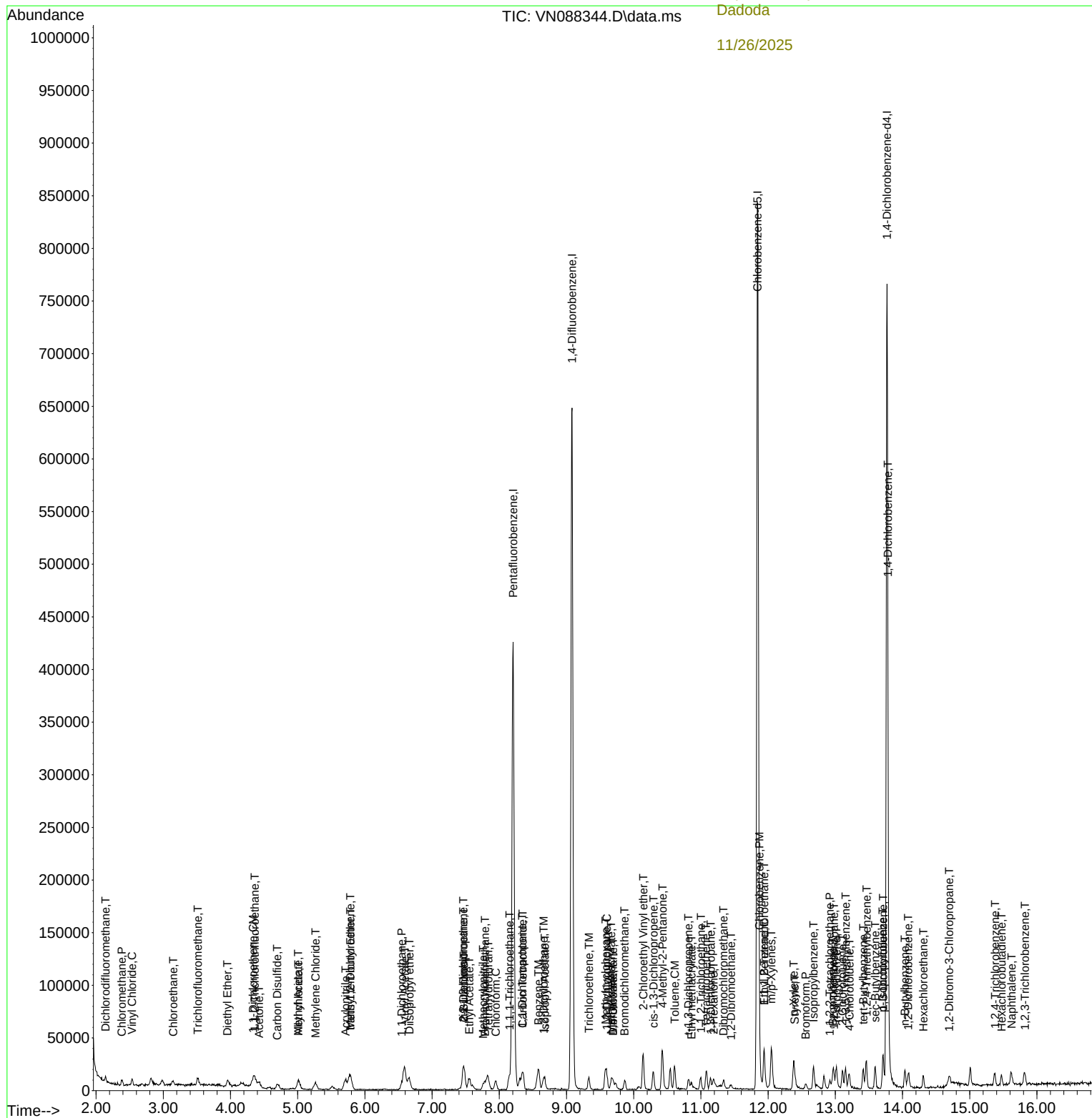
Manual Integrations APPROVED

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 : 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

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

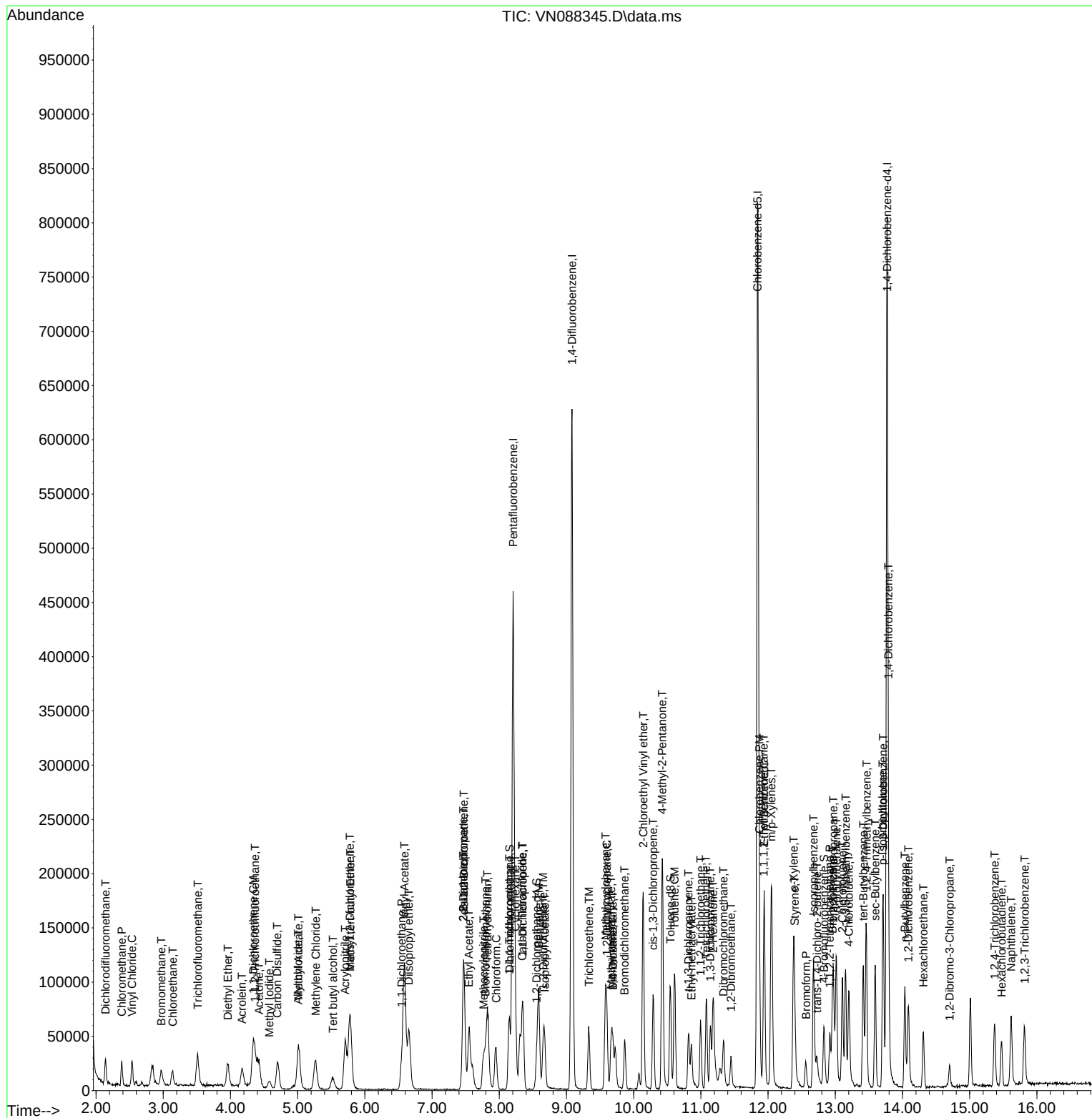
Quant Time: Nov 25 01:10:10 2025

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

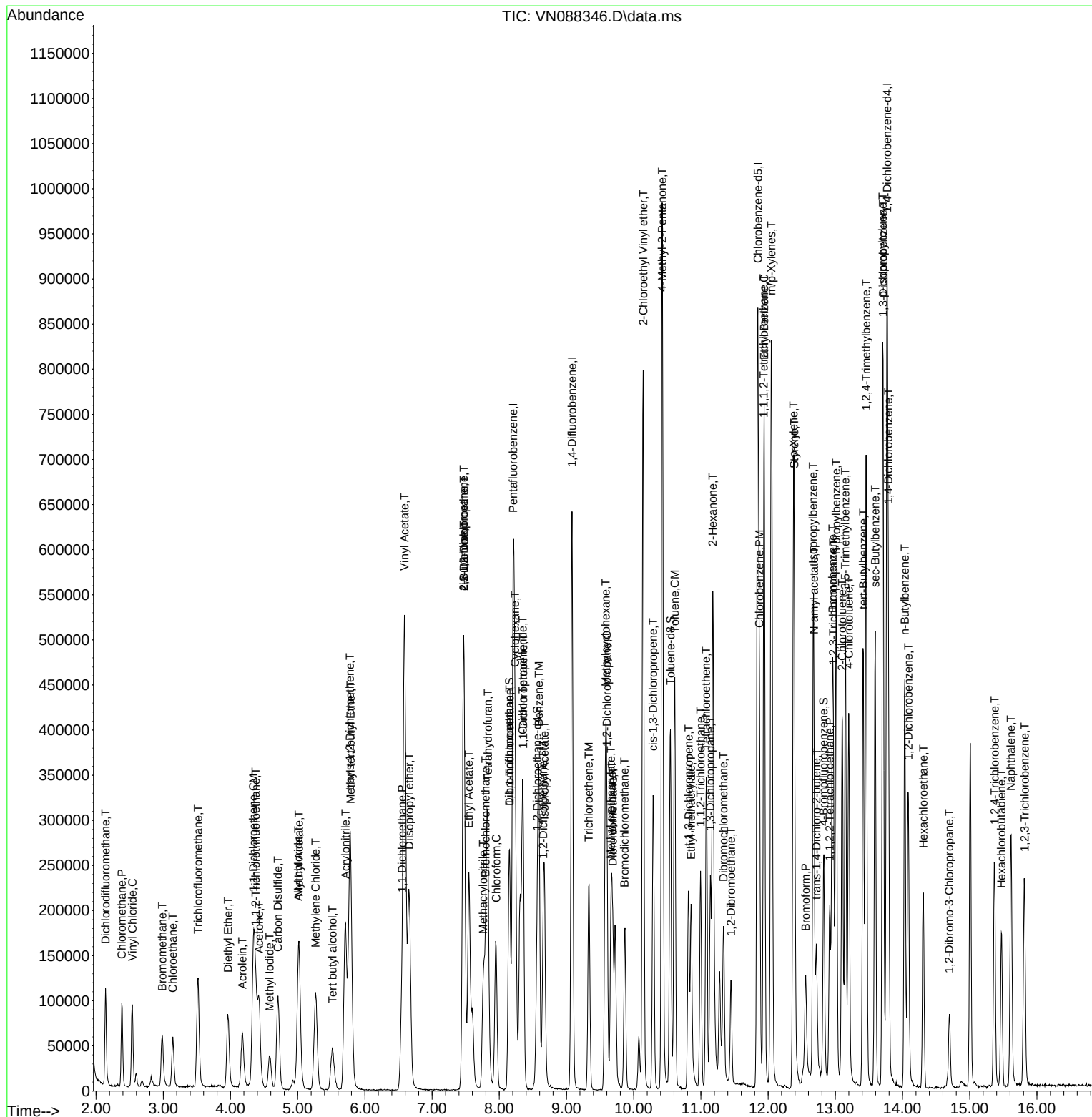
Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
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(#) = qualifier out of range (m) = manual integration (+) = signals summed

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

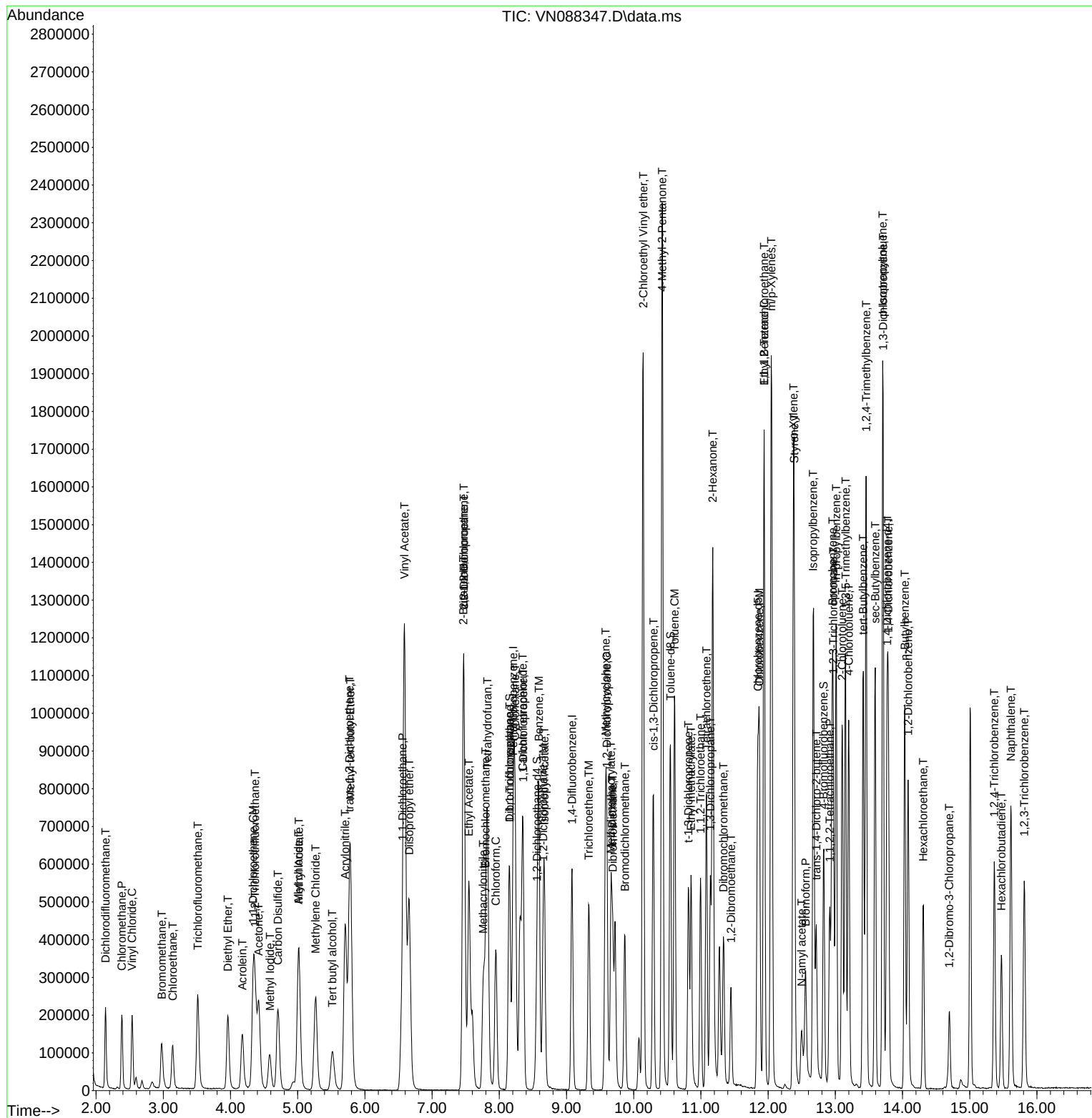
Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN112425\
Data File : VN088347.D
Acq On : 24 Nov 2025 13:53
Operator : JC\MD
Sample : VSTDICCC050
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 5 Sample Multiplier: 1

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

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)

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

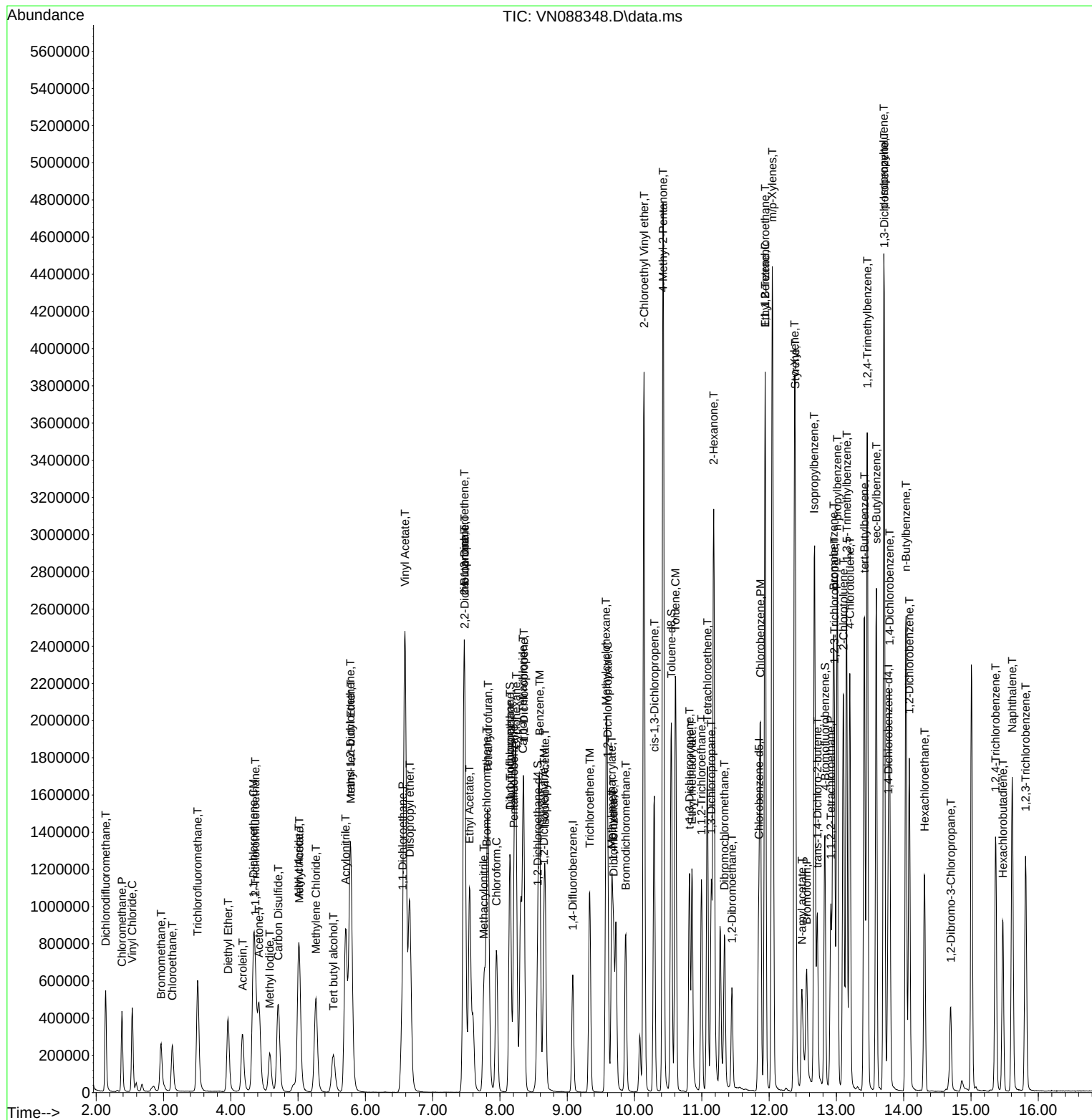
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Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN112425\  
Data File : VN088348.D  
Acq On    : 24 Nov 2025  14:14  
Operator  : JC\MD  
Sample    : VSTDICC100  
Misc      : 5.0mL/MSVOA_N/WATER  
ALS Vial  : 6    Sample Multiplier: 1
```

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



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

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

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

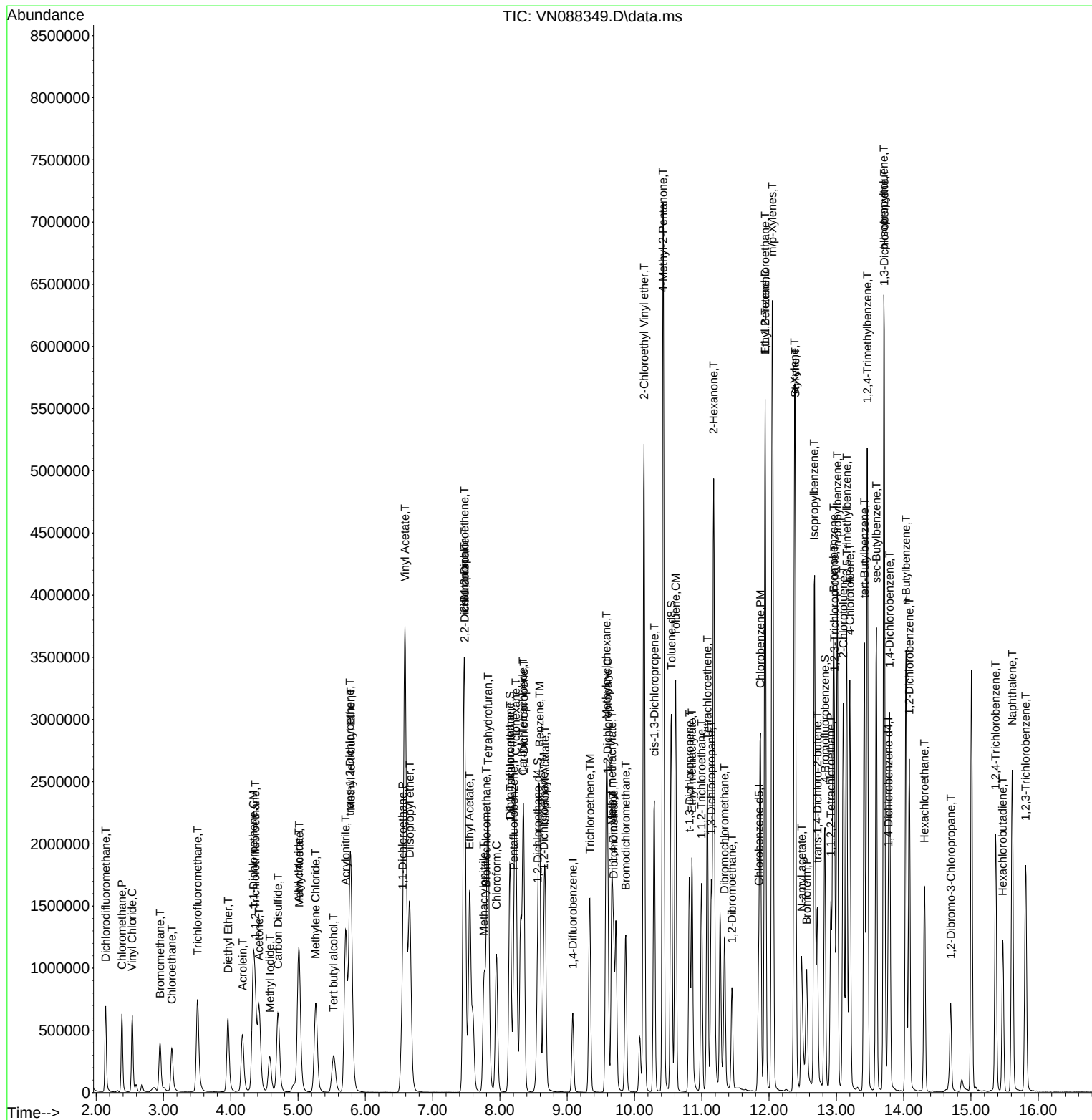
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Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN112425\  
Data File : VN088349.D  
Acq On    : 24 Nov 2025  14:35  
Operator  : JC\MD  
Sample    : VSTDICC150  
Misc      : 5.0mL/MSVOA_N/WATER  
ALS Vial  : 7    Sample Multiplier: 1
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Instrument :
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)
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(#) = 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-amyl 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 CONTINUING CALIBRATION CHECK

Lab Name:	<u>Alliance</u>	Contract:	<u>REMI01</u>
Lab Code:	<u>ACE</u>	SDG No.:	<u>Q3743</u>
Instrument ID:	<u>MSVOA_N</u>	Calibration Date/Time:	<u>12/01/2025 09:58</u>
Lab File ID:	<u>VN088370.D</u>	Init. Calib. Date(s):	<u>11/24/2025 11/24/2025</u>
Heated Purge: (Y/N)	<u>N</u>	Init. Calib. Time(s):	<u>12:38 14:35</u>
GC Column:	<u>RXI-624</u>	ID:	<u>0.25</u> (mm)

COMPOUND	RRF	RRF050	MIN RRF	%D	MAX%D
Dichlorodifluoromethane	0.751	0.815		8.52	20
Chloromethane	0.807	0.858	0.1	6.32	20
Vinyl Chloride	0.821	0.851		3.65	20
Bromomethane	0.395	0.408		3.29	20
Chloroethane	0.508	0.525		3.35	20
Trichlorofluoromethane	1.153	1.195		3.64	20
1,1,2-Trichlorotrifluoroethane	0.614	0.636		3.58	20
1,1-Dichloroethene	0.610	0.607		-0.49	20
Acetone	0.341	0.269		-21.11	20
Carbon Disulfide	1.993	2.083		4.52	20
Methyl tert-butyl Ether	2.285	2.626		14.92	20
Methyl Acetate	1.014	1.108		9.27	20
Methylene Chloride	0.729	0.791		8.51	20
trans-1,2-Dichloroethene	0.676	0.724		7.1	20
1,1-Dichloroethane	1.367	1.456	0.1	6.51	20
Cyclohexane	1.239	1.285		3.71	20
2-Butanone	0.548	0.541		-1.28	20
Carbon Tetrachloride	0.532	0.584		9.77	20
cis-1,2-Dichloroethene	0.785	0.861		9.68	20
Bromochloromethane	0.681	0.672		-1.32	20
Chloroform	1.343	1.443		7.45	20
1,1,1-Trichloroethane	1.181	1.220		3.3	20
Methylcyclohexane	0.572	0.664		16.08	20
Benzene	1.570	1.738		10.7	20
1,2-Dichloroethane	0.599	0.676		12.85	20
Trichloroethene	0.371	0.415		11.86	20
1,2-Dichloropropane	0.402	0.453		12.69	20
Bromodichloromethane	0.586	0.670		14.33	20
4-Methyl-2-Pentanone	0.633	0.747		18.01	20
Toluene	0.906	1.034		14.13	20
t-1,3-Dichloropropene	0.601	0.701		16.64	20
cis-1,3-Dichloropropene	0.632	0.748		18.35	20
1,1,2-Trichloroethane	0.351	0.400		13.96	20
2-Hexanone	0.399	0.458		14.79	20
Dibromochloromethane	0.400	0.464		16	20
1,2-Dibromoethane	0.355	0.413		16.34	20
Tetrachloroethene	0.396	0.425		7.32	20
Chlorobenzene	1.123	1.242	0.3	10.6	20

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



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>REMI01</u>
Lab Code:	<u>ACE</u>	SDG No.:	<u>Q3743</u>
Instrument ID:	<u>MSVOA_N</u>	Calibration Date/Time:	<u>12/01/2025</u> <u>09:58</u>
Lab File ID:	<u>VN088370.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
Ethyl Benzene	1.963	2.187		11.41	20
m/p-Xylenes	0.728	0.816		12.09	20
o-Xylene	0.694	0.778		12.1	20
Styrene	1.127	1.336		18.55	20
Bromoform	0.277	0.337	0.1	21.66	20
Isopropylbenzene	3.697	4.179		13.04	20
1,1,2,2-Tetrachloroethane	1.195	1.313	0.3	9.87	20
1,3-Dichlorobenzene	1.647	1.852		12.45	20
1,4-Dichlorobenzene	1.725	1.901		10.2	20
1,2-Dichlorobenzene	1.550	1.757		13.35	20
1,2-Dibromo-3-Chloropropane	0.290	0.317		9.31	20
1,2,4-Trichlorobenzene	0.921	1.042		13.14	20
1,2,3-Trichlorobenzene	0.896	0.989		10.38	20
1,2-Dichloroethane-d4	0.942	0.958		1.7	20
Dibromofluoromethane	0.346	0.377		8.96	20
Toluene-d8	1.289	1.356		5.2	20
4-Bromofluorobenzene	0.442	0.472		6.79	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\VN120125\
 Data File : VN088370.D
 Acq On : 01 Dec 2025 09:58
 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 12/02/2025
 Supervised By : Mahesh Dadoda 12/02/2025

Quant Time: Dec 02 00:07:59 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	8.200	168	262500	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.082	114	457607	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.841	117	408098	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.765	152	197460	50.000	ug/l	0.00

System Monitoring Compounds

33) 1,2-Dichloroethane-d4	8.559	65	251592	50.878	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	= 101.760%	
35) Dibromofluoromethane	8.147	113	172647	54.442	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 108.880%	
50) Toluene-d8	10.541	98	620606	52.597	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	= 105.200%	
62) 4-Bromofluorobenzene	12.823	95	216098	53.376	ug/l	0.00
Spiked Amount	50.000	Range	77 - 121	Recovery	= 106.760%	

Target Compounds

						Qvalue
2) Dichlorodifluoromethane	2.142	85	213988	54.274	ug/l	98
3) Chloromethane	2.383	50	225109	53.159	ug/l	97
4) Vinyl Chloride	2.536	62	223310	51.837	ug/l	98
5) Bromomethane	2.977	94	107123	51.665	ug/l	97
6) Chloroethane	3.136	64	137709	51.654	ug/l	96
7) Trichlorofluoromethane	3.512	101	313594	51.793	ug/l	100
8) Diethyl Ether	3.965	74	122135	53.743	ug/l	100
9) 1,1,2-Trichlorotrifluo...	4.365	101	166917	51.784	ug/l	97
10) Methyl Iodide	4.577	142	178922	56.072	ug/l	100
11) Tert butyl alcohol	5.518	59	205497	260.289	ug/l	97
12) 1,1-Dichloroethene	4.342	96	159359	49.760	ug/l	98
13) Acrolein	4.171	56	291964	386.274	ug/l	98
14) Allyl chloride	5.012	41	306169	51.579	ug/l	99
15) Acrylonitrile	5.706	53	631123	277.070	ug/l	99
16) Acetone	4.418	43	353017	197.397	ug/l	99
17) Carbon Disulfide	4.700	76	546678	52.238	ug/l	100
18) Methyl Acetate	5.012	43	290775	54.622	ug/l	100
19) Methyl tert-butyl Ether	5.783	73	689442	57.460	ug/l	96
20) Methylene Chloride	5.265	84	207762	54.299	ug/l	98
21) trans-1,2-Dichloroethene	5.771	96	190100	53.585	ug/l	98
22) Diisopropyl ether	6.653	45	704752	55.612	ug/l	98
23) Vinyl Acetate	6.589	43	2933303m	287.290	ug/l	
24) 1,1-Dichloroethane	6.553	63	382166	53.252	ug/l	99
25) 2-Butanone	7.465	43	709877	246.913	ug/l	99
26) 2,2-Dichloropropane	7.471	77	325483	52.543	ug/l	98
27) cis-1,2-Dichloroethene	7.471	96	225942	54.801	ug/l	98
28) Bromochloromethane	7.794	49	176525	49.377	ug/l	99
29) Tetrahydrofuran	7.824	42	580681	281.590	ug/l	99
30) Chloroform	7.947	83	378896	53.730	ug/l	95
31) Cyclohexane	8.235	56	337193	51.848	ug/l	98
32) 1,1,1-Trichloroethane	8.147	97	320304	51.659	ug/l	99
36) 1,1-Dichloropropene	8.353	75	257172	52.546	ug/l	98
37) Ethyl Acetate	7.541	43	351134	54.564	ug/l	98
38) Carbon Tetrachloride	8.341	117	267192	54.841	ug/l	95
39) Methylcyclohexane	9.582	83	303843	58.041	ug/l	97
40) Benzene	8.588	78	795502	55.380	ug/l	99

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN120125\
 Data File : VN088370.D
 Acq On : 01 Dec 2025 09:58
 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 12/02/2025

Supervised By :Mahesh Dadoda 12/02/2025

Quant Time: Dec 02 00:07:59 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	7.759	41	178298m	55.940	ug/l	
42) 1,2-Dichloroethane	8.653	62	309394	56.455	ug/l	99
43) Isopropyl Acetate	8.671	43	566144	57.580	ug/l	99
44) Trichloroethene	9.330	130	189717	55.896	ug/l	99
45) 1,2-Dichloropropane	9.600	63	207193	56.298	ug/l	94
46) Dibromomethane	9.688	93	152439	58.039	ug/l	98
47) Bromodichloromethane	9.865	83	306641	57.219	ug/l	99
48) Methyl methacrylate	9.659	41	262260	58.460	ug/l	99
49) 1,4-Dioxane	9.682	88	61264	1073.826	ug/l #	98
51) 4-Methyl-2-Pentanone	10.424	43	1709516	294.976	ug/l	99
52) Toluene	10.606	92	472964	57.023	ug/l	98
53) t-1,3-Dichloropropene	10.812	75	320705	58.302	ug/l	99
54) cis-1,3-Dichloropropene	10.294	75	342246	59.163	ug/l	96
55) 1,1,2-Trichloroethane	10.994	97	183113	57.061	ug/l	99
56) Ethyl methacrylate	10.853	69	317004	62.537	ug/l	98
57) 1,3-Dichloropropane	11.141	76	338648	58.496	ug/l	100
58) 2-Chloroethyl Vinyl ether	10.135	63	944366	323.148	ug/l	99
59) 2-Hexanone	11.176	43	1048679	287.241	ug/l	99
60) Dibromochloromethane	11.335	129	212408	58.062	ug/l	98
61) 1,2-Dibromoethane	11.447	107	189124	58.199	ug/l	99
64) Tetrachloroethene	11.082	164	173317	53.571	ug/l	99
65) Chlorobenzene	11.871	112	506678	55.274	ug/l	98
66) 1,1,1,2-Tetrachloroethane	11.935	131	176540	57.102	ug/l	100
67) Ethyl Benzene	11.941	91	892546	55.709	ug/l	99
68) m/p-Xylenes	12.047	106	666195	112.150	ug/l	99
69) o-Xylene	12.376	106	317531	56.069	ug/l	98
70) Styrene	12.388	104	545300	59.301	ug/l	99
71) Bromoform	12.559	173	137488	60.715	ug/l #	100
73) Isopropylbenzene	12.670	105	825111	56.515	ug/l	99
74) N-amyl acetate	12.494	43	246831m	49.643	ug/l	
75) 1,1,2,2-Tetrachloroethane	12.918	83	259360	54.950	ug/l	99
76) 1,2,3-Trichloropropane	12.970	75	258773m	55.953	ug/l	
77) Bromobenzene	12.959	156	190979	57.171	ug/l	98
78) n-propylbenzene	13.012	91	1008099	56.441	ug/l	100
79) 2-Chlorotoluene	13.100	91	601237	55.123	ug/l	99
80) 1,3,5-Trimethylbenzene	13.147	105	692333	57.227	ug/l	100
81) trans-1,4-Dichloro-2-b...	12.718	75	97180	59.611	ug/l #	87
82) 4-Chlorotoluene	13.200	91	633461	58.473	ug/l	99
83) tert-Butylbenzene	13.412	119	573074	55.723	ug/l	100
84) 1,2,4-Trimethylbenzene	13.459	105	704771	58.609	ug/l	99
85) sec-Butylbenzene	13.594	105	840504	56.889	ug/l	100
86) p-Isopropyltoluene	13.706	119	696965	57.528	ug/l	99
87) 1,3-Dichlorobenzene	13.712	146	365660	56.231	ug/l	99
88) 1,4-Dichlorobenzene	13.788	146	375462	55.121	ug/l	99
89) n-Butylbenzene	14.029	91	663659	56.418	ug/l	100
90) Hexachloroethane	14.306	117	124701	56.344	ug/l	99
91) 1,2-Dichlorobenzene	14.082	146	346992	56.704	ug/l	99
92) 1,2-Dibromo-3-Chloropr...	14.694	75	62609	54.672	ug/l	98
93) 1,2,4-Trichlorobenzene	15.364	180	205849	56.615	ug/l	98
94) Hexachlorobutadiene	15.470	225	69826	53.442	ug/l	98
95) Naphthalene	15.611	128	732311	58.123	ug/l	99
96) 1,2,3-Trichlorobenzene	15.811	180	195344	55.205	ug/l	97

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN120125\
Data File : VN088370.D
Acq On : 01 Dec 2025 09:58
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 12/02/2025
Supervised By :Mahesh Dadoda 12/02/2025

Quant Time: Dec 02 00:07:59 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)
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(#) = qualifier out of range (m) = manual integration (+) = signals summed

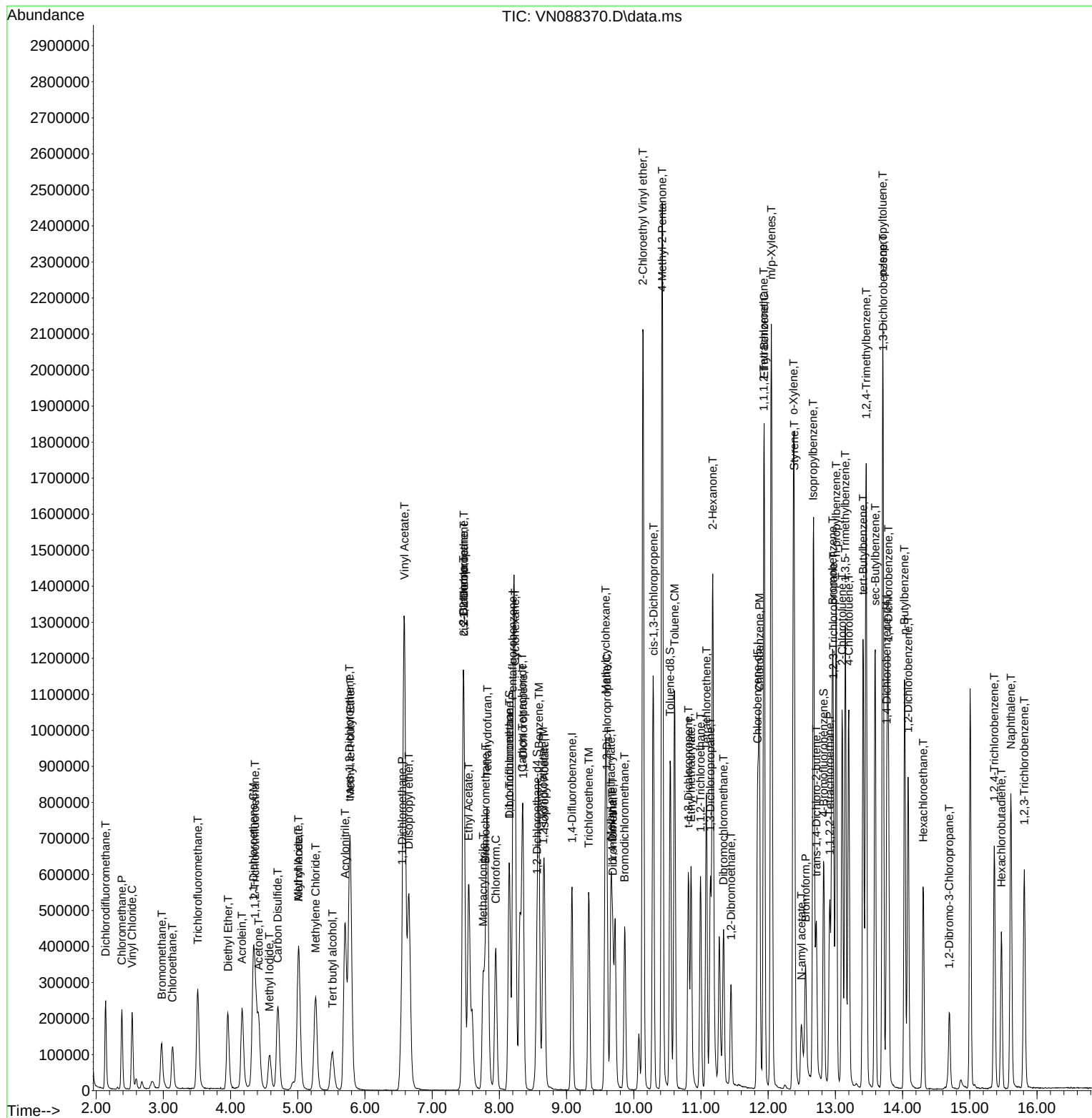
Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN120125\
Data File : VN088370.D
Acq On : 01 Dec 2025 09:58
Operator : JC\MD
Sample : VSTDCCC050
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 3 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
VSTDCCC050

Manual Integrations APPROVED

Reviewed By :John Carlone 12/02/2025
Supervised By :Mahesh Dadoda 12/02/2025

Quant Time: Dec 02 00:07:59 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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN120125\
 Data File : VN088370.D
 Acq On : 01 Dec 2025 09:58
 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 02 00:07:59 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

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	103	0.00
2 T	Dichlorodifluoromethane	0.751	0.815	-8.5	114	0.00
3 P	Chloromethane	0.807	0.858	-6.3	110	0.00
4 C	Vinyl Chloride	0.821	0.851	-3.7#	108	0.00
5 T	Bromomethane	0.395	0.408	-3.3	106	0.00
6 T	Chloroethane	0.508	0.525	-3.3	107	0.00
7 T	Trichlorofluoromethane	1.153	1.195	-3.6	113	0.00
8 T	Diethyl Ether	0.433	0.465	-7.4	108	0.00
9 T	1,1,2-Trichlorotrifluoroeth	0.614	0.636	-3.6	116	0.00
10 T	Methyl Iodide	0.608	0.682	-12.2	106	-0.01
11 T	Tert butyl alcohol	0.150	0.157	-4.7	100	0.00
12 CM	1,1-Dichloroethene	0.610	0.607	0.5#	110	0.00
13 T	Acrolein	0.144	0.222	-54.2#	152#	0.00
14 T	Allyl chloride	1.131	1.166	-3.1	105	0.00
15 T	Acrylonitrile	0.434	0.481	-10.8	106	0.00
16 T	Acetone	0.341	0.269	21.1	83	0.00
17 T	Carbon Disulfide	1.993	2.083	-4.5	108	0.00
18 T	Methyl Acetate	1.014	1.108	-9.3	109	0.00
19 T	Methyl tert-butyl Ether	2.285	2.626	-14.9	108	0.00
20 T	Methylene Chloride	0.729	0.791	-8.5	108	0.00
21 T	trans-1,2-Dichloroethene	0.676	0.724	-7.1	109	0.00
22 T	Diisopropyl ether	2.414	2.685	-11.2	107	0.00
23 T	Vinyl Acetate	1.945	2.235	-14.9	108	0.00
24 P	1,1-Dichloroethane	1.367	1.456	-6.5	105	0.00
25 T	2-Butanone	0.548	0.541	1.3	94	0.00
26 T	2,2-Dichloropropane	1.180	1.240	-5.1	106	0.00
27 T	cis-1,2-Dichloroethene	0.785	0.861	-9.7	107	0.00
28 T	Bromochloromethane	0.681	0.672	1.3	94	0.00
29 T	Tetrahydrofuran	0.393	0.442	-12.5	106	0.00
30 C	Chloroform	1.343	1.443	-7.4#	107	0.00
31 T	Cyclohexane	1.239	1.285	-3.7	117	0.00
32 T	1,1,1-Trichloroethane	1.181	1.220	-3.3	109	0.00
33 S	1,2-Dichloroethane-d4	0.942	0.958	-1.7	99	0.00
34 I	1,4-Difluorobenzene	1.000	1.000	0.0	95	0.00
35 S	Dibromofluoromethane	0.346	0.377	-9.0	103	0.00
36 T	1,1-Dichloropropene	0.535	0.562	-5.0	109	0.00
37 T	Ethyl Acetate	0.703	0.767	-9.1	101	0.00
38 T	Carbon Tetrachloride	0.532	0.584	-9.8	111	0.00
39 T	Methylcyclohexane	0.572	0.664	-16.1	120	0.00
40 TM	Benzene	1.570	1.738	-10.7	106	0.00
41 T	Methacrylonitrile	0.348	0.390	-12.1	104	0.00
42 TM	1,2-Dichloroethane	0.599	0.676	-12.9	105	0.00
43 T	Isopropyl Acetate	1.074	1.237	-15.2	105	0.00
44 TM	Trichloroethene	0.371	0.415	-11.9	110	0.00
45 C	1,2-Dichloropropane	0.402	0.453	-12.7#	106	0.00
46 T	Dibromomethane	0.287	0.333	-16.0	108	0.00
47 T	Bromodichloromethane	0.586	0.670	-14.3	108	0.00
48 T	Methyl methacrylate	0.490	0.573	-16.9	106	0.00

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN120125\
 Data File : VN088370.D
 Acq On : 01 Dec 2025 09:58
 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 02 00:07:59 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

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	0.007	-16.7	96	0.00
50 S	Toluene-d8	1.289	1.356	-5.2	100	0.00
51 T	4-Methyl-2-Pentanone	0.633	0.747	-18.0	106	0.00
52 CM	Toluene	0.906	1.034	-14.1#	108	0.00
53 T	t-1,3-Dichloropropene	0.601	0.701	-16.6	106	0.00
54 T	cis-1,3-Dichloropropene	0.632	0.748	-18.4	108	0.00
55 T	1,1,2-Trichloroethane	0.351	0.400	-14.0	108	0.00
56 T	Ethyl methacrylate	0.554	0.693	-25.1#	107	0.00
57 T	1,3-Dichloropropane	0.633	0.740	-16.9	109	0.00
58 T	2-Chloroethyl Vinyl ether	0.319	0.413	-29.5#	108	0.00
59 T	2-Hexanone	0.399	0.458	-14.8	100	0.00
60 T	Dibromochloromethane	0.400	0.464	-16.0	108	0.00
61 T	1,2-Dibromoethane	0.355	0.413	-16.3	106	0.00
62 S	4-Bromofluorobenzene	0.442	0.472	-6.8	100	0.00
63 I	Chlorobenzene-d5	1.000	1.000	0.0	96	0.00
64 T	Tetrachloroethene	0.396	0.425	-7.3	108	0.00
65 PM	Chlorobenzene	1.123	1.242	-10.6	108	0.00
66 T	1,1,1,2-Tetrachloroethane	0.379	0.433	-14.2	108	0.00
67 C	Ethyl Benzene	1.963	2.187	-11.4#	108	0.00
68 T	m/p-Xylenes	0.728	0.816	-12.1	107	0.00
69 T	o-Xylene	0.694	0.778	-12.1	107	0.00
70 T	Styrene	1.127	1.336	-18.5	107	0.00
71 P	Bromoform	0.277	0.337	-21.7	113	0.00
72 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	95	0.00
73 T	Isopropylbenzene	3.697	4.179	-13.0	109	0.00
74 T	N-amyl acetate	1.259	1.250	0.7	113	0.00
75 P	1,1,2,2-Tetrachloroethane	1.195	1.313	-9.9	110	0.00
76 T	1,2,3-Trichloropropane	1.171	1.311	-12.0	103	0.00
77 T	Bromobenzene	0.846	0.967	-14.3	106	0.00
78 T	n-propylbenzene	4.523	5.105	-12.9	109	0.00
79 T	2-Chlorotoluene	2.762	3.045	-10.2	107	0.00
80 T	1,3,5-Trimethylbenzene	3.063	3.506	-14.5	109	0.00
81 T	trans-1,4-Dichloro-2-butene	0.413	0.492	-19.1	108	0.00
82 T	4-Chlorotoluene	2.743	3.208	-17.0	106	0.00
83 T	tert-Butylbenzene	2.604	2.902	-11.4	109	0.00
84 T	1,2,4-Trimethylbenzene	3.045	3.569	-17.2	109	0.00
85 T	sec-Butylbenzene	3.741	4.257	-13.8	112	0.00
86 T	p-Isopropyltoluene	3.068	3.530	-15.1	110	0.00
87 T	1,3-Dichlorobenzene	1.647	1.852	-12.4	109	0.00
88 T	1,4-Dichlorobenzene	1.725	1.901	-10.2	107	0.00
89 T	n-Butylbenzene	2.979	3.361	-12.8	111	0.00
90 T	Hexachloroethane	0.560	0.632	-12.9	112	0.00
91 T	1,2-Dichlorobenzene	1.550	1.757	-13.4	109	0.00
92 T	1,2-Dibromo-3-Chloropropane	0.290	0.317	-9.3	104	0.00
93 T	1,2,4-Trichlorobenzene	0.921	1.042	-13.1	111	0.00
94 T	Hexachlorobutadiene	0.331	0.354	-6.9	113	0.00
95 T	Naphthalene	3.190	3.709	-16.3	107	0.00

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN120125\
Data File : VN088370.D
Acq On : 01 Dec 2025 09:58
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 02 00:07:59 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

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	0.989	-10.4	107	0.00

(#) = Out of Range

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN120125\
 Data File : VN088370.D
 Acq On : 01 Dec 2025 09:58
 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 02 00:07:59 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

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	103	0.00
2 T	Dichlorodifluoromethane	50.000	54.274	-8.5	114	0.00
3 P	Chloromethane	50.000	53.159	-6.3	110	0.00
4 C	Vinyl Chloride	50.000	51.837	-3.7#	108	0.00
5 T	Bromomethane	50.000	51.665	-3.3	106	0.00
6 T	Chloroethane	50.000	51.654	-3.3	107	0.00
7 T	Trichlorofluoromethane	50.000	51.793	-3.6	113	0.00
8 T	Diethyl Ether	50.000	53.743	-7.5	108	0.00
9 T	1,1,2-Trichlorotrifluoroeth	50.000	51.784	-3.6	116	0.00
10 T	Methyl Iodide	50.000	56.072	-12.1	106	-0.01
11 T	Tert butyl alcohol	250.000	260.289	-4.1	100	0.00
12 CM	1,1-Dichloroethene	50.000	49.760	0.5#	110	0.00
13 T	Acrolein	250.000	386.274	-54.5#	152	0.00
14 T	Allyl chloride	50.000	51.579	-3.2	105	0.00
15 T	Acrylonitrile	250.000	277.070	-10.8	106	0.00
16 T	Acetone	250.000	197.397	21.0	83	0.00
17 T	Carbon Disulfide	50.000	52.238	-4.5	108	0.00
18 T	Methyl Acetate	50.000	54.622	-9.2	109	0.00
19 T	Methyl tert-butyl Ether	50.000	57.460	-14.9	108	0.00
20 T	Methylene Chloride	50.000	54.299	-8.6	108	0.00
21 T	trans-1,2-Dichloroethene	50.000	53.585	-7.2	109	0.00
22 T	Diisopropyl ether	50.000	55.612	-11.2	107	0.00
23 T	Vinyl Acetate	250.000	287.290	-14.9	108	0.00
24 P	1,1-Dichloroethane	50.000	53.252	-6.5	105	0.00
25 T	2-Butanone	250.000	246.913	1.2	94	0.00
26 T	2,2-Dichloropropane	50.000	52.543	-5.1	106	0.00
27 T	cis-1,2-Dichloroethene	50.000	54.801	-9.6	107	0.00
28 T	Bromochloromethane	50.000	49.377	1.2	94	0.00
29 T	Tetrahydrofuran	250.000	281.590	-12.6	106	0.00
30 C	Chloroform	50.000	53.730	-7.5#	107	0.00
31 T	Cyclohexane	50.000	51.848	-3.7	117	0.00
32 T	1,1,1-Trichloroethane	50.000	51.659	-3.3	109	0.00
33 S	1,2-Dichloroethane-d4	50.000	50.878	-1.8	99	0.00
34 I	1,4-Difluorobenzene	50.000	50.000	0.0	95	0.00
35 S	Dibromofluoromethane	50.000	54.442	-8.9	103	0.00
36 T	1,1-Dichloropropene	50.000	52.546	-5.1	109	0.00
37 T	Ethyl Acetate	50.000	54.564	-9.1	101	0.00
38 T	Carbon Tetrachloride	50.000	54.841	-9.7	111	0.00
39 T	Methylcyclohexane	50.000	58.041	-16.1	120	0.00
40 TM	Benzene	50.000	55.380	-10.8	106	0.00
41 T	Methacrylonitrile	50.000	55.940	-11.9	104	0.00
42 TM	1,2-Dichloroethane	50.000	56.455	-12.9	105	0.00
43 T	Isopropyl Acetate	50.000	57.580	-15.2	105	0.00
44 TM	Trichloroethene	50.000	55.896	-11.8	110	0.00
45 C	1,2-Dichloropropane	50.000	56.298	-12.6#	106	0.00
46 T	Dibromomethane	50.000	58.039	-16.1	108	0.00
47 T	Bromodichloromethane	50.000	57.219	-14.4	108	0.00
48 T	Methyl methacrylate	50.000	58.460	-16.9	106	0.00

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN120125\
 Data File : VN088370.D
 Acq On : 01 Dec 2025 09:58
 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 02 00:07:59 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

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	1073.826	-7.4	96	0.00
50 S	Toluene-d8	50.000	52.597	-5.2	100	0.00
51 T	4-Methyl-2-Pentanone	250.000	294.976	-18.0	106	0.00
52 CM	Toluene	50.000	57.023	-14.0#	108	0.00
53 T	t-1,3-Dichloropropene	50.000	58.302	-16.6	106	0.00
54 T	cis-1,3-Dichloropropene	50.000	59.163	-18.3	108	0.00
55 T	1,1,2-Trichloroethane	50.000	57.061	-14.1	108	0.00
56 T	Ethyl methacrylate	50.000	62.537	-25.1#	107	0.00
57 T	1,3-Dichloropropane	50.000	58.496	-17.0	109	0.00
58 T	2-Chloroethyl Vinyl ether	250.000	323.148	-29.3#	108	0.00
59 T	2-Hexanone	250.000	287.241	-14.9	100	0.00
60 T	Dibromochloromethane	50.000	58.062	-16.1	108	0.00
61 T	1,2-Dibromoethane	50.000	58.199	-16.4	106	0.00
62 S	4-Bromofluorobenzene	50.000	53.376	-6.8	100	0.00
63 I	Chlorobenzene-d5	50.000	50.000	0.0	96	0.00
64 T	Tetrachloroethene	50.000	53.571	-7.1	108	0.00
65 PM	Chlorobenzene	50.000	55.274	-10.5	108	0.00
66 T	1,1,1,2-Tetrachloroethane	50.000	57.102	-14.2	108	0.00
67 C	Ethyl Benzene	50.000	55.709	-11.4#	108	0.00
68 T	m/p-Xylenes	100.000	112.150	-12.2	107	0.00
69 T	o-Xylene	50.000	56.069	-12.1	107	0.00
70 T	Styrene	50.000	59.301	-18.6	107	0.00
71 P	Bromoform	50.000	60.715	-21.4	113	0.00
72 I	1,4-Dichlorobenzene-d4	50.000	50.000	0.0	95	0.00
73 T	Isopropylbenzene	50.000	56.515	-13.0	109	0.00
74 T	N-amyl acetate	50.000	49.643	0.7	113	0.00
75 P	1,1,2,2-Tetrachloroethane	50.000	54.950	-9.9	110	0.00
76 T	1,2,3-Trichloropropane	50.000	55.953	-11.9	103	0.00
77 T	Bromobenzene	50.000	57.171	-14.3	106	0.00
78 T	n-propylbenzene	50.000	56.441	-12.9	109	0.00
79 T	2-Chlorotoluene	50.000	55.123	-10.2	107	0.00
80 T	1,3,5-Trimethylbenzene	50.000	57.227	-14.5	109	0.00
81 T	trans-1,4-Dichloro-2-butene	50.000	59.611	-19.2	108	0.00
82 T	4-Chlorotoluene	50.000	58.473	-16.9	106	0.00
83 T	tert-Butylbenzene	50.000	55.723	-11.4	109	0.00
84 T	1,2,4-Trimethylbenzene	50.000	58.609	-17.2	109	0.00
85 T	sec-Butylbenzene	50.000	56.889	-13.8	112	0.00
86 T	p-Isopropyltoluene	50.000	57.528	-15.1	110	0.00
87 T	1,3-Dichlorobenzene	50.000	56.231	-12.5	109	0.00
88 T	1,4-Dichlorobenzene	50.000	55.121	-10.2	107	0.00
89 T	n-Butylbenzene	50.000	56.418	-12.8	111	0.00
90 T	Hexachloroethane	50.000	56.344	-12.7	112	0.00
91 T	1,2-Dichlorobenzene	50.000	56.704	-13.4	109	0.00
92 T	1,2-Dibromo-3-Chloropropane	50.000	54.672	-9.3	104	0.00
93 T	1,2,4-Trichlorobenzene	50.000	56.615	-13.2	111	0.00
94 T	Hexachlorobutadiene	50.000	53.442	-6.9	113	0.00
95 T	Naphthalene	50.000	58.123	-16.2	107	0.00

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN120125\
Data File : VN088370.D
Acq On : 01 Dec 2025 09:58
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 02 00:07:59 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

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	55.205	-10.4	107	0.00

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



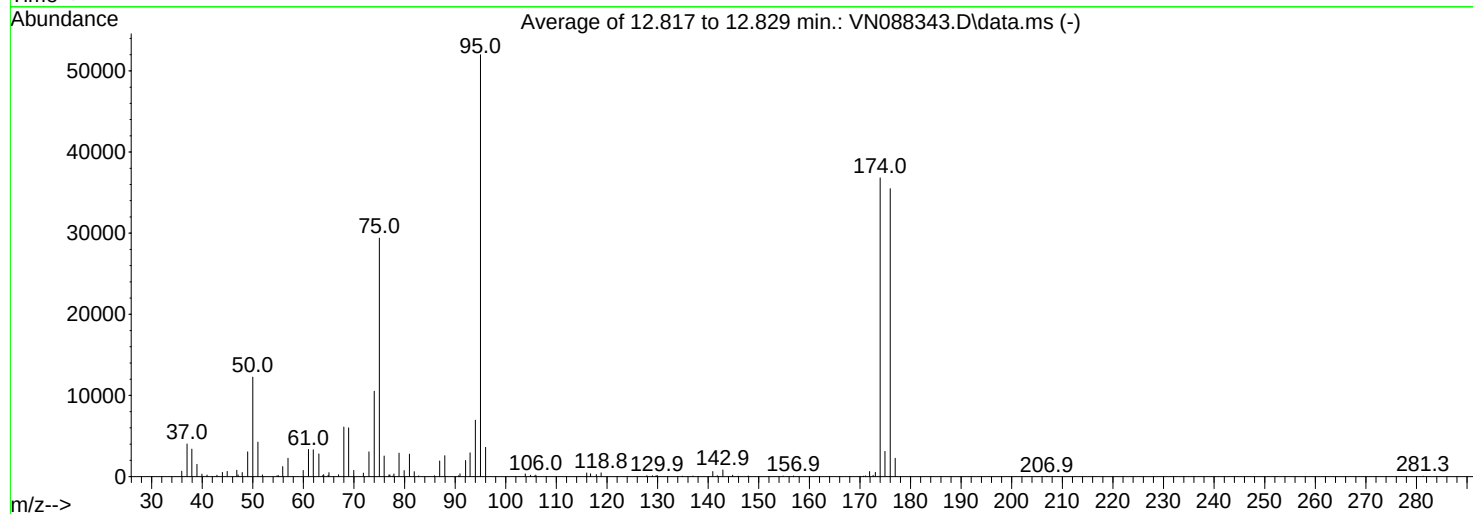
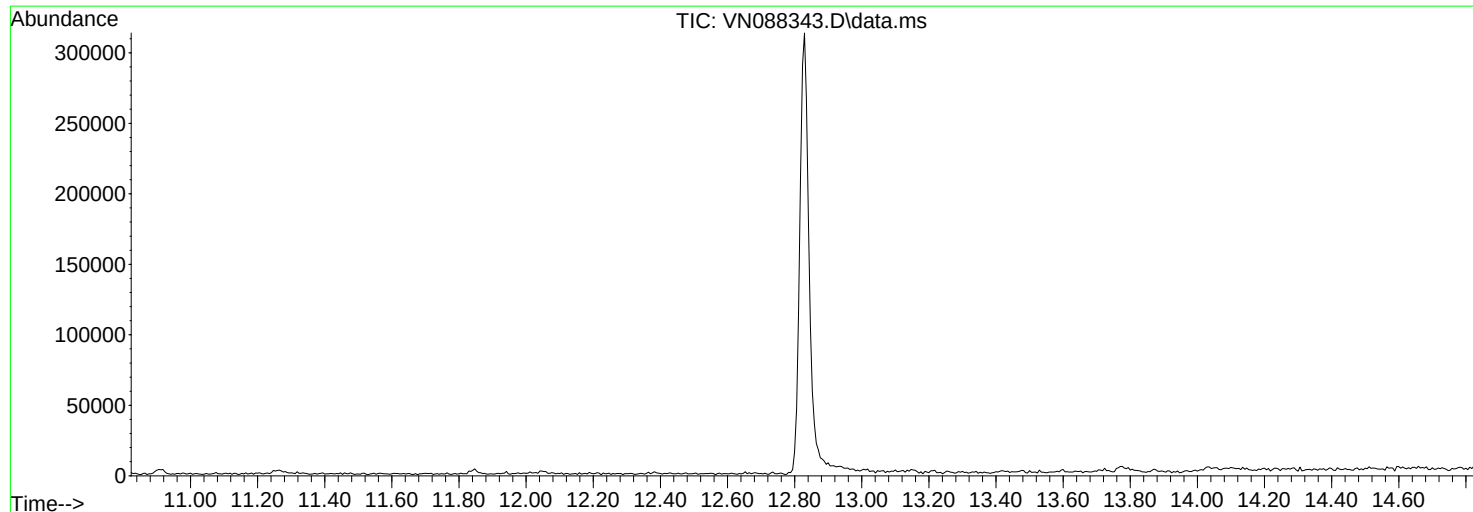
QC SAMPLE DATA

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

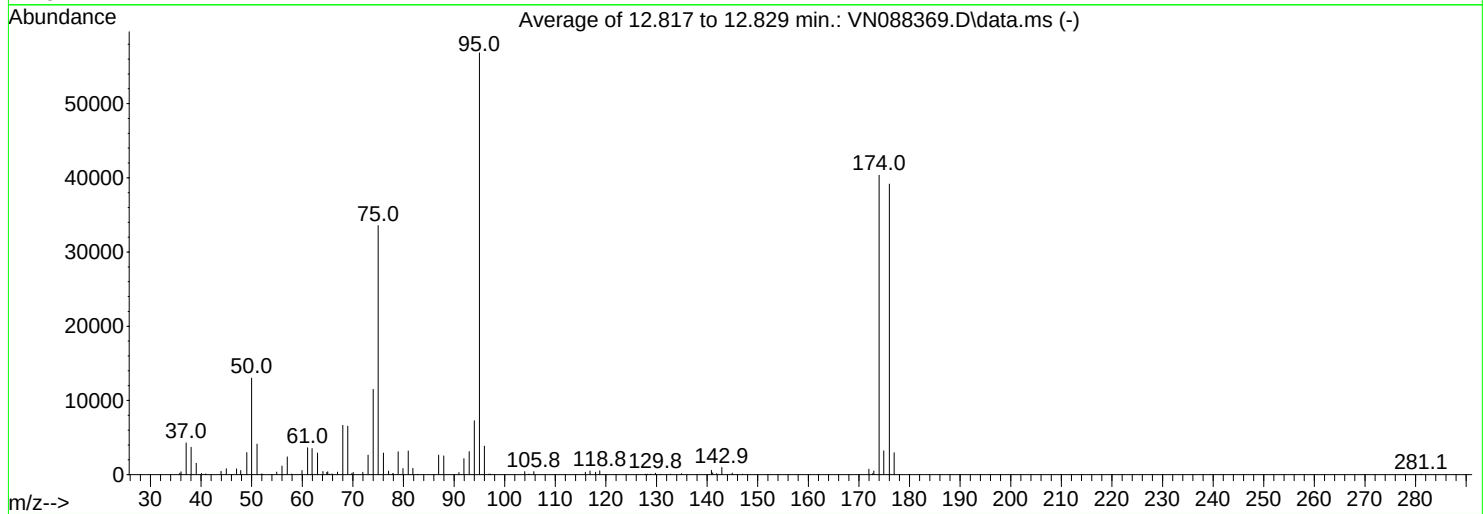
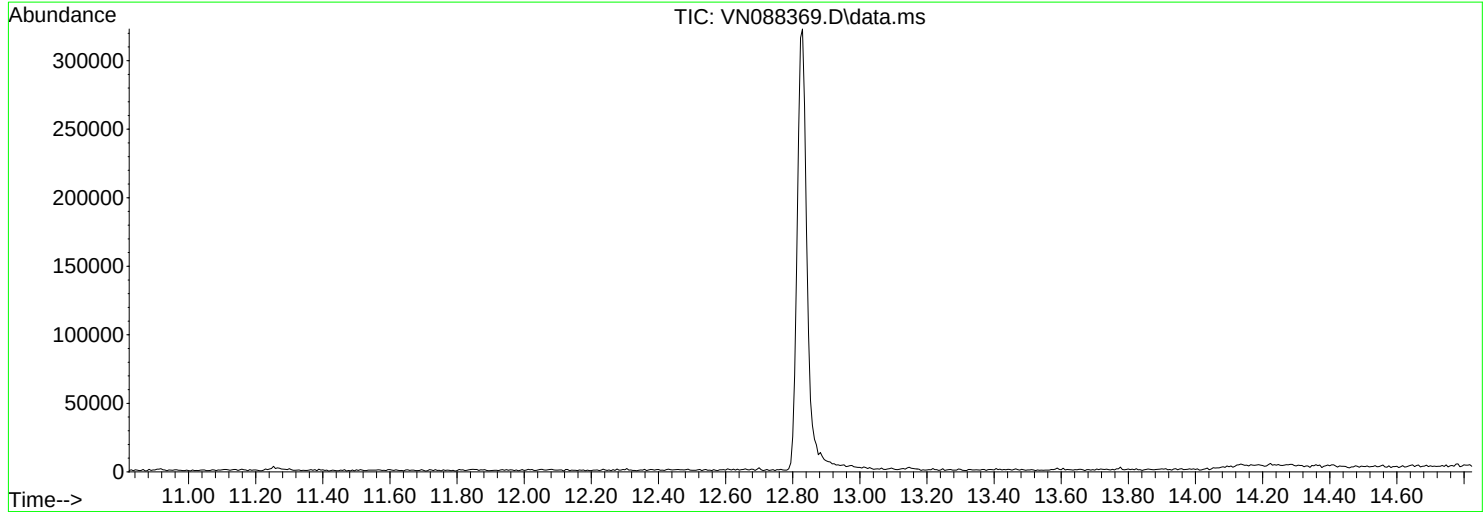
m/z	abund.	m/z	abund.	m/z	abund.	m/z	abund.
281.30	53						

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN120125\
Data File : VN088369.D
Acq On : 01 Dec 2025 08:49
Operator : JC\MD
Sample : BFB
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 2 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 1838

Target Mass	Rel. to Mass	Lower Limit%	Upper Limit%	Rel. Abn%	Raw Abn	Result Pass/Fail
50	95	15	40	22.9	13006	PASS
75	95	30	60	59.1	33576	PASS
95	95	100	100	100.0	56845	PASS
96	95	5	9	6.7	3835	PASS
173	174	0.00	2	1.2	489	PASS
174	95	50	100	71.0	40341	PASS
175	174	5	9	8.0	3213	PASS
176	174	95	101	97.1	39165	PASS
177	176	5	9	7.5	2949	PASS

BFB

Modified:subtracted

m/z	abund.	m/z	abund.	m/z	abund.	m/z	abund.
35.80	190	49.05	2976	61.95	3517	70.05	311
36.05	384	50.00	13006	63.00	2917	71.95	331
37.05	4272	51.05	4113	64.10	433	73.00	263
38.05	3684	51.90	88	64.85	304	74.00	1149
39.05	1525	52.10	68	65.05	399	75.00	33576
40.10	142	54.95	339	65.40	72	76.05	2917
40.95	105	56.00	1161	65.90	102	77.05	468
43.95	455	57.05	2400	66.95	355	77.80	128
45.00	808	57.90	81	68.00	6639	78.00	126
47.00	782	59.95	560	69.00	6543	78.95	3079
47.85	567	61.05	3625	69.80	174	79.90	824

Average of 12.817 to 12.829 min.: VN088369.D\data.ms

BFB

Modified:subtracted

m/z	abund.	m/z	abund.	m/z	abund.	m/z	abund.
80.95	3198	96.00	3835	128.20	67	147.00	61
81.85	841	96.90	89	128.80	61	148.00	50
82.80	51	97.20	58	129.80	188	171.95	743
86.00	50	103.95	408	130.80	50	172.70	124
86.95	2627	104.80	82	132.80	67	172.95	489
87.95	2534	105.75	414	134.95	138	174.00	40341
90.95	286	115.95	332	140.85	577	174.90	3213
91.95	2167	116.85	505	141.10	238	176.00	39165
93.00	3106	117.95	315	141.90	151	176.95	2949
94.00	7272	118.80	513	142.90	969	281.10	73
95.00	56845	127.80	56	144.95	206		

Report of Analysis

Client: Remington & Vernick
 Project: Saddler Property
 Client Sample ID: VN1201WBL01
 Lab Sample ID: VN1201WBL01
 Analytical Method: 8260D
 Sample Wt/Vol: 5 mL

Level: LOW
 Final Vol: 5000 uL

Date Collected:
 Date Received:
 SDG No.: Q3743
 Matrix: Water
 % Solid: 0
 Test: VOC-TCLVOA-10

CAS Number	Parameter	Conc.	Qua.	DF	MDL	LOQ / CRQL	Units	Date Ana.	BatchID
TARGETS									
75-71-8	Dichlorodifluoromethane	0.22	U	1	0.22	1.00	ug/L	12/01/25 10:53	VN120125
74-87-3	Chloromethane	0.32	U	1	0.32	1.00	ug/L	12/01/25 10:53	VN120125
75-01-4	Vinyl Chloride	0.26	U	1	0.26	1.00	ug/L	12/01/25 10:53	VN120125
74-83-9	Bromomethane	1.40	U	1	1.40	5.00	ug/L	12/01/25 10:53	VN120125
75-00-3	Chloroethane	0.47	U	1	0.47	1.00	ug/L	12/01/25 10:53	VN120125
75-69-4	Trichlorofluoromethane	0.33	U	1	0.33	1.00	ug/L	12/01/25 10:53	VN120125
76-13-1	1,1,2-Trichlorotrifluoroethane	0.25	U	1	0.25	1.00	ug/L	12/01/25 10:53	VN120125
75-35-4	1,1-Dichloroethene	0.23	U	1	0.23	1.00	ug/L	12/01/25 10:53	VN120125
67-64-1	Acetone	1.50	U	1	1.50	5.00	ug/L	12/01/25 10:53	VN120125
75-15-0	Carbon Disulfide	0.21	U	1	0.21	1.00	ug/L	12/01/25 10:53	VN120125
1634-04-4	Methyl tert-butyl Ether	0.16	U	1	0.16	1.00	ug/L	12/01/25 10:53	VN120125
79-20-9	Methyl Acetate	0.27	U	1	0.27	1.00	ug/L	12/01/25 10:53	VN120125
75-09-2	Methylene Chloride	0.28	U	1	0.28	1.00	ug/L	12/01/25 10:53	VN120125
156-60-5	trans-1,2-Dichloroethene	0.23	U	1	0.23	1.00	ug/L	12/01/25 10:53	VN120125
75-34-3	1,1-Dichloroethane	0.23	U	1	0.23	1.00	ug/L	12/01/25 10:53	VN120125
110-82-7	Cyclohexane	1.50	U	1	1.50	5.00	ug/L	12/01/25 10:53	VN120125
78-93-3	2-Butanone	0.98	U	1	0.98	5.00	ug/L	12/01/25 10:53	VN120125
56-23-5	Carbon Tetrachloride	0.25	U	1	0.25	1.00	ug/L	12/01/25 10:53	VN120125
156-59-2	cis-1,2-Dichloroethene	0.19	U	1	0.19	1.00	ug/L	12/01/25 10:53	VN120125
74-97-5	Bromochloromethane	0.22	U	1	0.22	1.00	ug/L	12/01/25 10:53	VN120125
67-66-3	Chloroform	0.25	U	1	0.25	1.00	ug/L	12/01/25 10:53	VN120125
71-55-6	1,1,1-Trichloroethane	0.20	U	1	0.20	1.00	ug/L	12/01/25 10:53	VN120125
108-87-2	Methylcyclohexane	0.16	U	1	0.16	1.00	ug/L	12/01/25 10:53	VN120125
71-43-2	Benzene	0.15	U	1	0.15	1.00	ug/L	12/01/25 10:53	VN120125
107-06-2	1,2-Dichloroethane	0.22	U	1	0.22	1.00	ug/L	12/01/25 10:53	VN120125
79-01-6	Trichloroethene	0.090	U	1	0.090	1.00	ug/L	12/01/25 10:53	VN120125
78-87-5	1,2-Dichloropropane	0.20	U	1	0.20	1.00	ug/L	12/01/25 10:53	VN120125
75-27-4	Bromodichloromethane	0.22	U	1	0.22	1.00	ug/L	12/01/25 10:53	VN120125
108-10-1	4-Methyl-2-Pentanone	0.68	U	1	0.68	5.00	ug/L	12/01/25 10:53	VN120125
108-88-3	Toluene	0.14	U	1	0.14	1.00	ug/L	12/01/25 10:53	VN120125
10061-02-6	t-1,3-Dichloropropene	0.17	U	1	0.17	1.00	ug/L	12/01/25 10:53	VN120125
10061-01-5	cis-1,3-Dichloropropene	0.16	U	1	0.16	1.00	ug/L	12/01/25 10:53	VN120125
79-00-5	1,1,2-Trichloroethane	0.21	U	1	0.21	1.00	ug/L	12/01/25 10:53	VN120125
591-78-6	2-Hexanone	0.89	U	1	0.89	5.00	ug/L	12/01/25 10:53	VN120125
124-48-1	Dibromochloromethane	0.18	U	1	0.18	1.00	ug/L	12/01/25 10:53	VN120125
106-93-4	1,2-Dibromoethane	0.15	U	1	0.15	1.00	ug/L	12/01/25 10:53	VN120125
127-18-4	Tetrachloroethene	0.23	U	1	0.23	1.00	ug/L	12/01/25 10:53	VN120125
108-90-7	Chlorobenzene	0.12	U	1	0.12	1.00	ug/L	12/01/25 10:53	VN120125
100-41-4	Ethyl Benzene	0.13	U	1	0.13	1.00	ug/L	12/01/25 10:53	VN120125
179601-23-1	m/p-Xylenes	0.24	U	1	0.24	2.00	ug/L	12/01/25 10:53	VN120125
95-47-6	o-Xylene	0.12	U	1	0.12	1.00	ug/L	12/01/25 10:53	VN120125
100-42-5	Styrene	0.15	U	1	0.15	1.00	ug/L	12/01/25 10:53	VN120125
75-25-2	Bromoform	0.19	U	1	0.19	1.00	ug/L	12/01/25 10:53	VN120125



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

Report of Analysis

Client: Remington & Vernick
Project: Saddler Property
Client Sample ID: VN1201WBL01
Lab Sample ID: VN1201WBL01
Analytical Method: 8260D
Sample Wt/Vol: 5 mL

Level: LOW
Final Vol: 5000 uL

Date Collected:
Date Received:
SDG No.: Q3743
Matrix: Water
% Solid: 0
Test: VOC-TCLVOA-10

CAS Number	Parameter	Conc.	Qua.	DF	MDL	LOQ / CRQL	Units	Date Ana.	BatchID
98-82-8	Isopropylbenzene	0.12	U	1	0.12	1.00	ug/L	12/01/25 10:53	VN120125
79-34-5	1,1,2,2-Tetrachloroethane	0.26	U	1	0.26	1.00	ug/L	12/01/25 10:53	VN120125
541-73-1	1,3-Dichlorobenzene	0.16	U	1	0.16	1.00	ug/L	12/01/25 10:53	VN120125
106-46-7	1,4-Dichlorobenzene	0.19	U	1	0.19	1.00	ug/L	12/01/25 10:53	VN120125
95-50-1	1,2-Dichlorobenzene	0.16	U	1	0.16	1.00	ug/L	12/01/25 10:53	VN120125
96-12-8	1,2-Dibromo-3-Chloropropane	0.53	U	1	0.53	1.00	ug/L	12/01/25 10:53	VN120125
120-82-1	1,2,4-Trichlorobenzene	0.20	U	1	0.20	1.00	ug/L	12/01/25 10:53	VN120125
87-61-6	1,2,3-Trichlorobenzene	0.20	U	1	0.20	1.00	ug/L	12/01/25 10:53	VN120125

SURROGATES

17060-07-0	1,2-Dichloroethane-d4	57.8			74 - 125	116%	SPK: 50
1868-53-7	Dibromofluoromethane	52.2			75 - 124	104%	SPK: 50
2037-26-5	Toluene-d8	50.0			86 - 113	100%	SPK: 50
460-00-4	4-Bromofluorobenzene	52.4			77 - 121	105%	SPK: 50

INTERNAL STANDARDS

INTERNAL STANDARDS		Area Count
363-72-4	Pentafluorobenzene	188000
540-36-3	1,4-Difluorobenzene	435000
3114-55-4	Chlorobenzene-d5	418000
3855-82-1	1,4-Dichlorobenzene-d4	199000

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\VN120125\
 Data File : VN088372.D
 Acq On : 01 Dec 2025 10:53
 Operator : JC\MD
 Sample : VN1201WBL01
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 VN1201WBL01

Quant Time: Dec 02 00:09:38 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	8.200	168	188076	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.082	114	435319	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.841	117	418422	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.770	152	198799	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.559	65	204928	57.841	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	= 115.680%	
35) Dibromofluoromethane	8.147	113	157356	52.161	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 104.320%	
50) Toluene-d8	10.541	98	561082	49.987	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	= 99.980%	
62) 4-Bromofluorobenzene	12.829	95	201912	52.426	ug/l	0.00
Spiked Amount	50.000	Range	77 - 121	Recovery	= 104.860%	

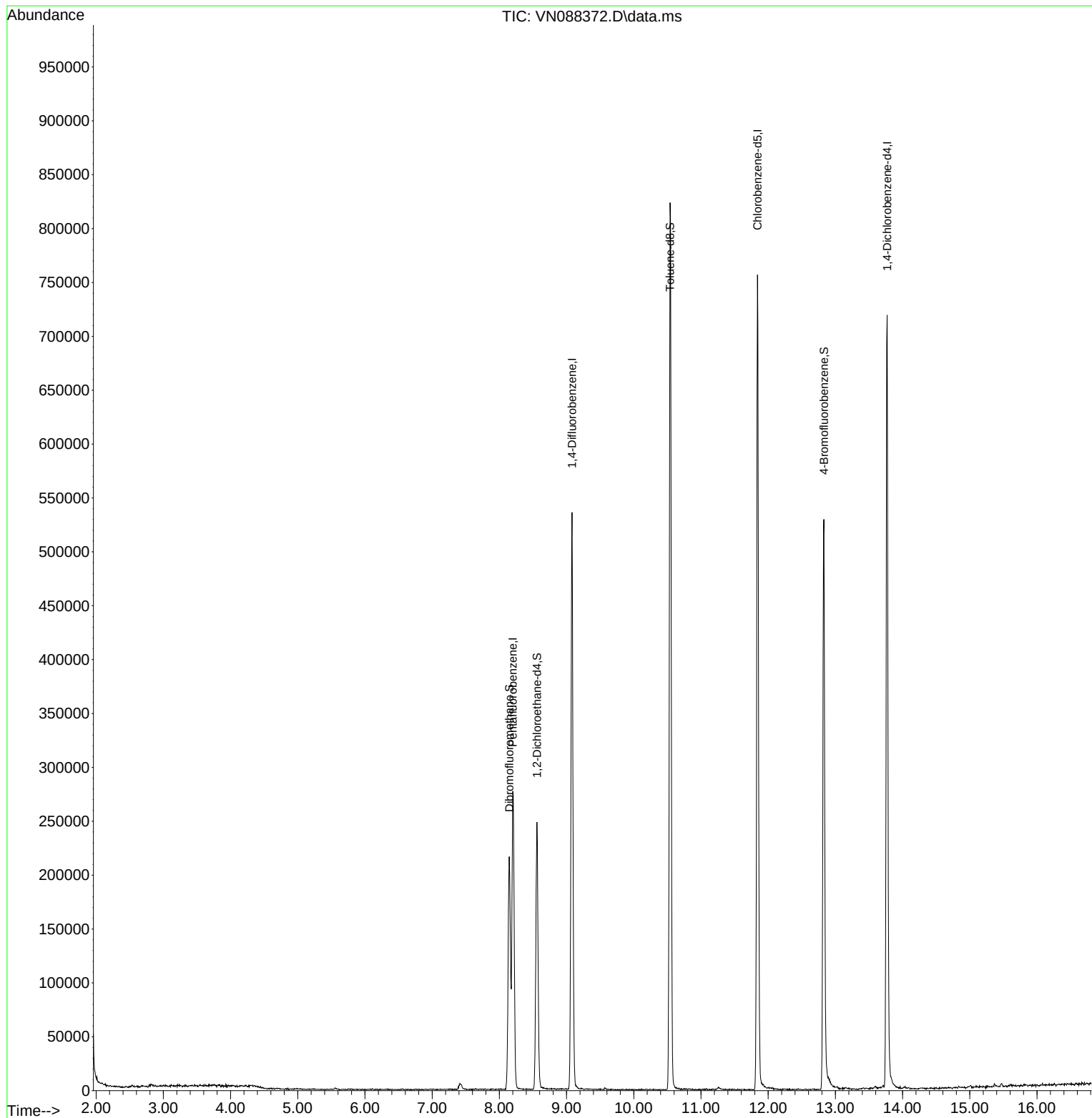
Target Compounds	Qvalue

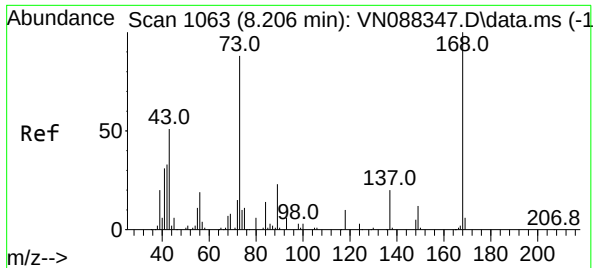
(#) = qualifier out of range (m) = manual integration (+) = signals summed

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN120125\
Data File : VN088372.D
Acq On : 01 Dec 2025 10:53
Operator : JC\MD
Sample : VN1201WBL01
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 5 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
VN1201WBL01

Quant Time: Dec 02 00:09:38 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

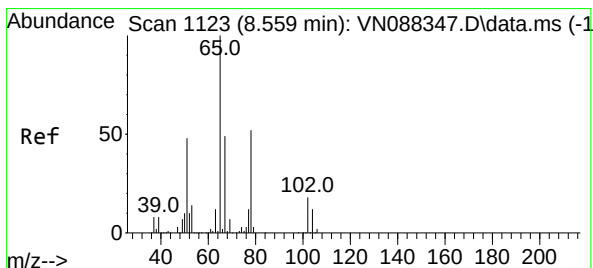
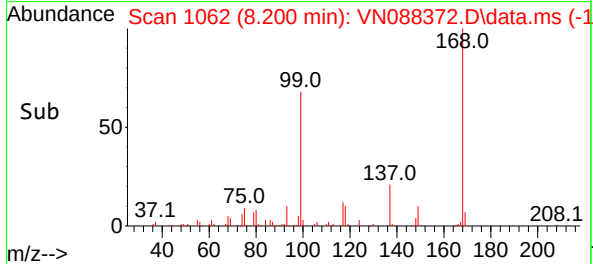
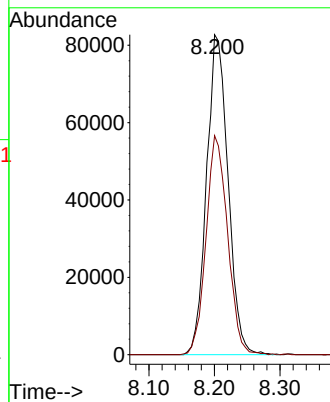
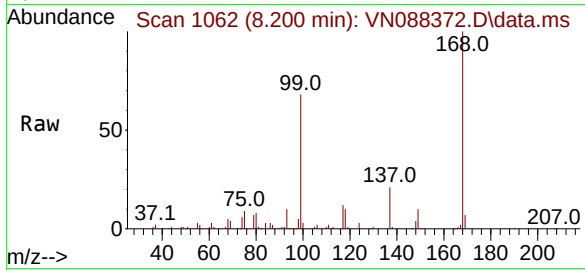




#1
Pentafluorobenzene
Concen: 50.000 ug/l
RT: 8.200 min Scan# 1063
Delta R.T. -0.006 min
Lab File: VN088372.D
Acq: 01 Dec 2025 10:53

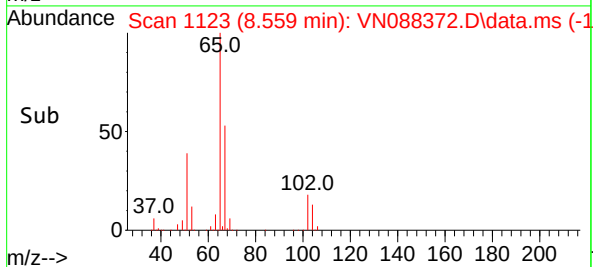
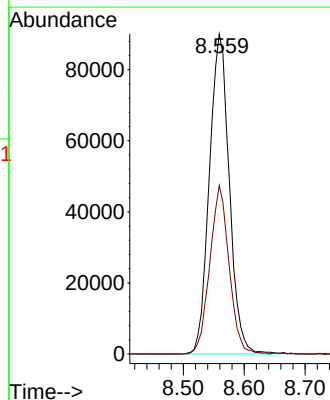
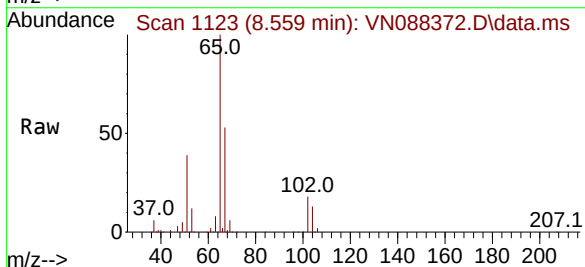
Instrument :
MSVOA_N
ClientSampleId :
VN1201WBL01

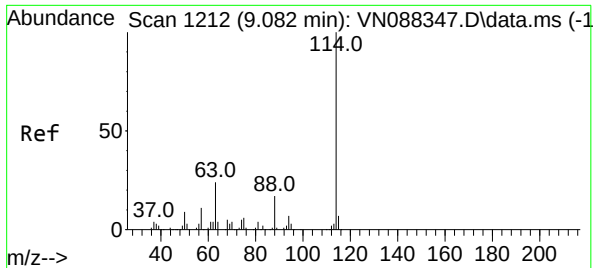
Tgt Ion:168 Resp: 188076
Ion Ratio Lower Upper
168 100
99 68.5 57.1 85.7



#33
1,2-Dichloroethane-d4
Concen: 57.841 ug/l
RT: 8.559 min Scan# 1123
Delta R.T. 0.000 min
Lab File: VN088372.D
Acq: 01 Dec 2025 10:53

Tgt Ion: 65 Resp: 204928
Ion Ratio Lower Upper
65 100
67 50.8 0.0 100.8





#34

1,4-Difluorobenzene

Concen: 50.000 ug/l

RT: 9.082 min Scan# 1

Delta R.T. 0.000 min

Lab File: VN088372.D

Acq: 01 Dec 2025 10:53

Instrument :

MSVOA_N

ClientSampleId :

VN1201WBL01

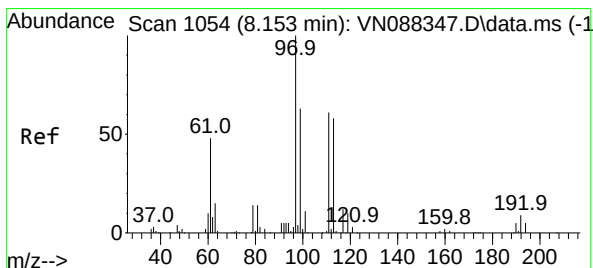
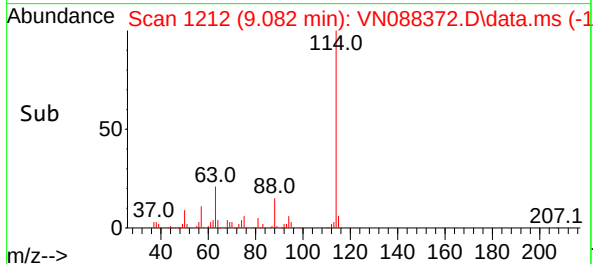
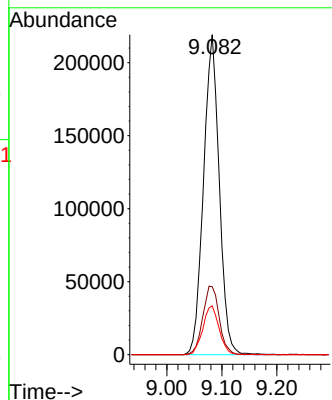
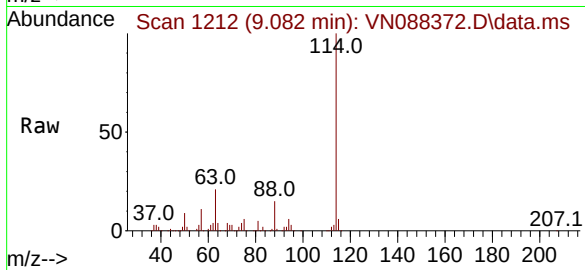
Tgt Ion:114 Resp: 435319

Ion Ratio Lower Upper

114 100

63 21.3 0.0 49.0

88 15.3 0.0 34.0



#35

Dibromofluoromethane

Concen: 52.161 ug/l

RT: 8.147 min Scan# 1053

Delta R.T. -0.006 min

Lab File: VN088372.D

Acq: 01 Dec 2025 10:53

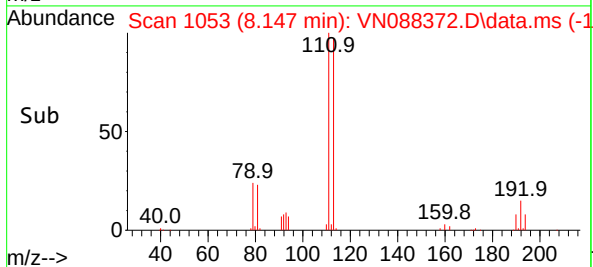
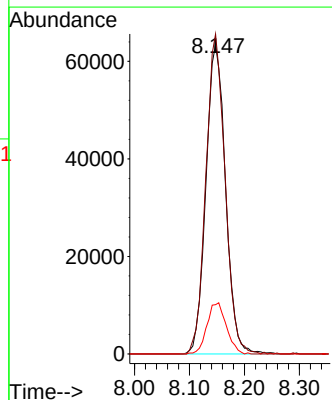
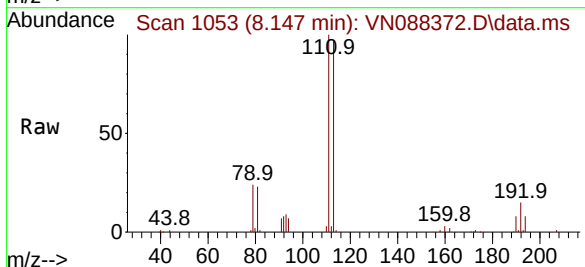
Tgt Ion:113 Resp: 157356

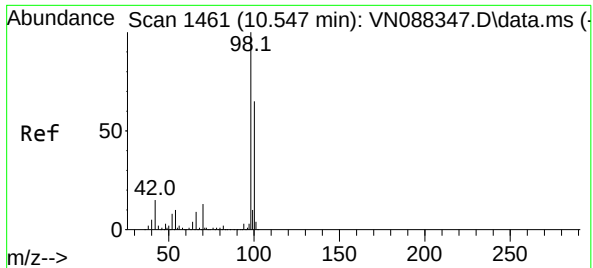
Ion Ratio Lower Upper

113 100

111 101.1 82.5 123.7

192 16.3 12.8 19.2

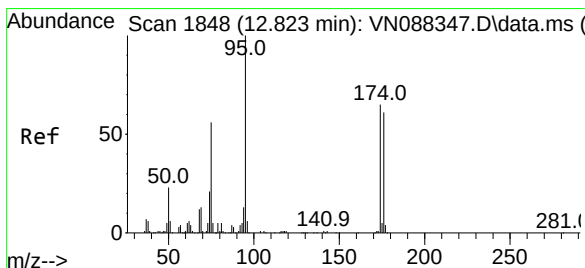
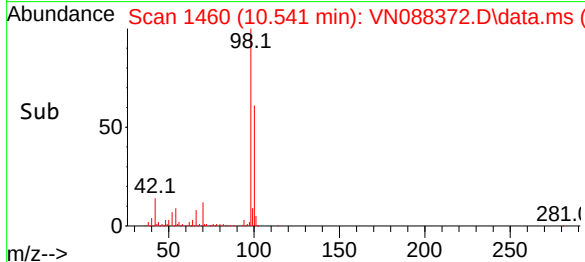
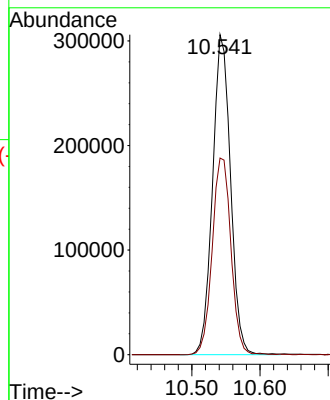
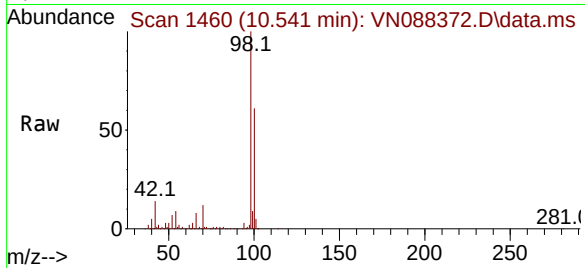




#50
Toluene-d8
Concen: 49.987 ug/l
RT: 10.541 min Scan# 1460
Delta R.T. -0.006 min
Lab File: VN088372.D
Acq: 01 Dec 2025 10:53

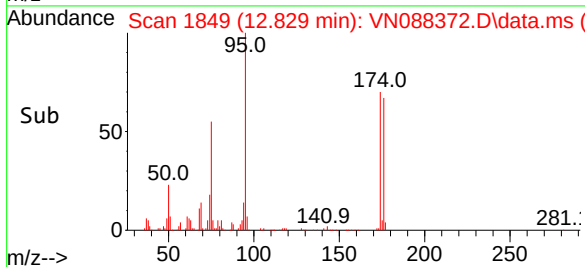
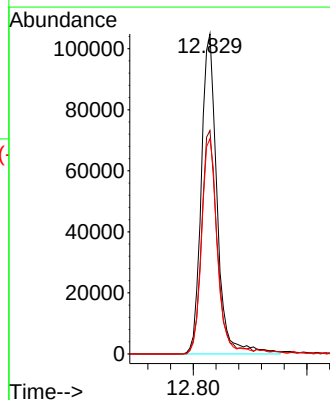
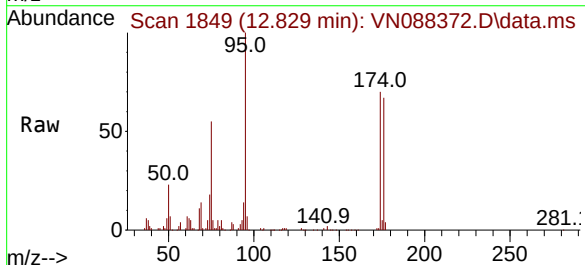
Instrument :
MSVOA_N
ClientSampleId :
VN1201WBL01

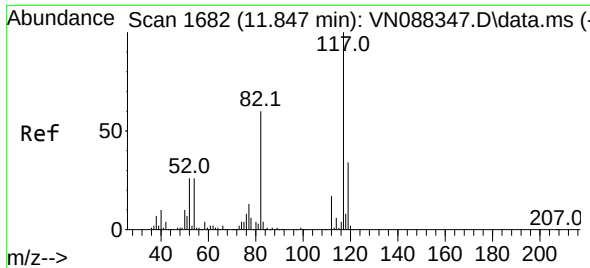
Tgt Ion: 98 Resp: 561082
Ion Ratio Lower Upper
98 100
100 64.0 52.2 78.2



#62
4-Bromofluorobenzene
Concen: 52.426 ug/l
RT: 12.829 min Scan# 1849
Delta R.T. 0.006 min
Lab File: VN088372.D
Acq: 01 Dec 2025 10:53

Tgt Ion: 95 Resp: 201912
Ion Ratio Lower Upper
95 100
174 69.2 0.0 139.0
176 67.0 0.0 132.4

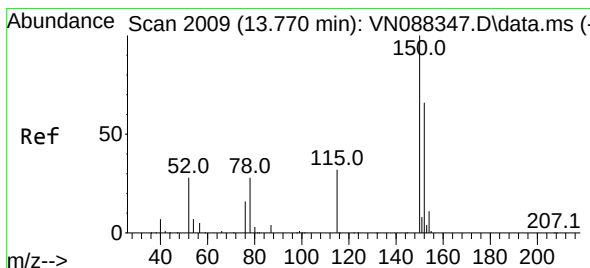
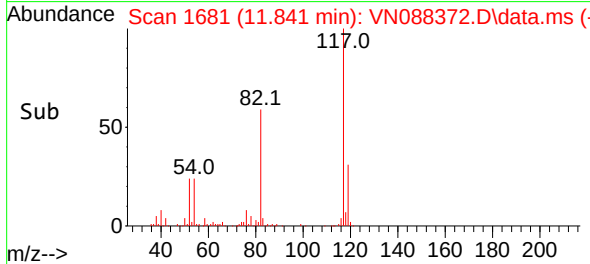
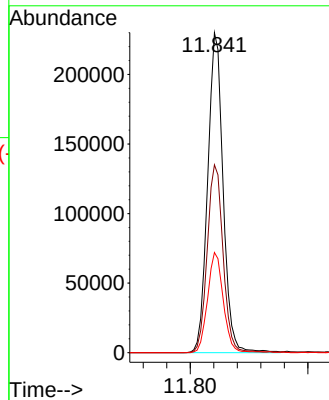
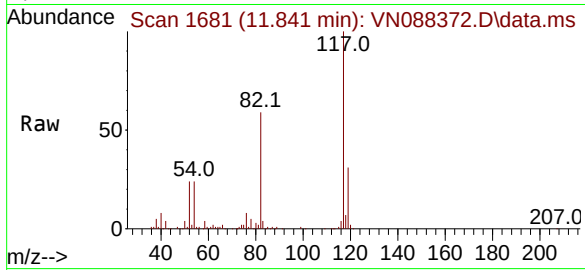




#63
Chlorobenzene-d5
Concen: 50.000 ug/l
RT: 11.841 min Scan# 11
Delta R.T. -0.006 min
Lab File: VN088372.D
Acq: 01 Dec 2025 10:53

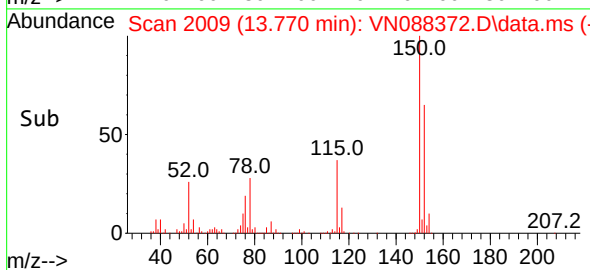
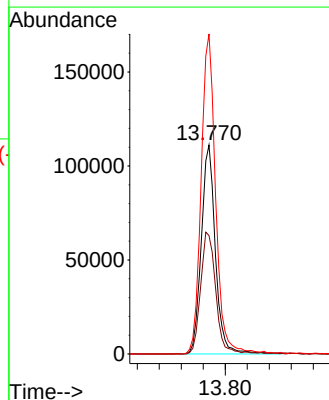
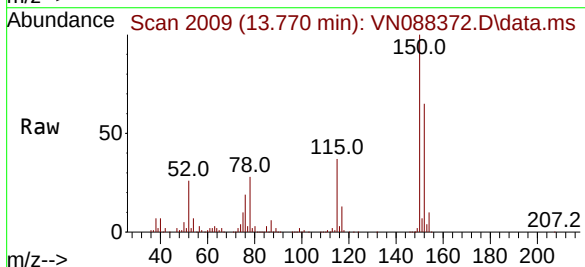
Instrument :
MSVOA_N
ClientSampleId :
VN1201WBL01

Tgt Ion	Ratio	Lower	Upper
117	100		
82	58.7	47.8	71.8
119	31.3	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: VN088372.D
Acq: 01 Dec 2025 10:53

Tgt Ion	Ratio	Lower	Upper
152	100		
115	62.5	33.0	98.9
150	156.8	0.0	349.0



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN120125\
Data File : VN088372.D
Acq On : 01 Dec 2025 10:53
Operator : JC\MD
Sample : VN1201WBL01
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 5 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
VN1201WBL01

Integration Parameters: RTEINT.P

Integrator: RTE

Smoothing : ON

Sampling : 1

Start Thrs: 0.2

Stop Thrs : 0

Filtering: 5

Min Area: 3 % of largest Peak

Max Peaks: 100

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

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

Title : SW846 8260

Signal : TIC: VN088372.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	7.412	921	928	937	rBV3	5682	16399	1.06%	0.202%
2	8.147	1042	1053	1058	rBV2	216362	540527	34.79%	6.646%
3	8.200	1058	1062	1077	rVB	276320	627960	40.42%	7.721%
4	8.559	1114	1123	1133	rBV	247884	561166	36.12%	6.900%
5	9.082	1203	1212	1228	rBV	535181	1099272	70.76%	13.516%
6	10.541	1452	1460	1475	rBV	822937	1553554	100.00%	19.101%
7	11.841	1672	1681	1694	rBV	756205	1399467	90.08%	17.207%
8	12.829	1841	1849	1859	rBV	529059	1004686	64.67%	12.353%
9	13.770	2001	2009	2027	rBV	716850	1330258	85.63%	16.356%

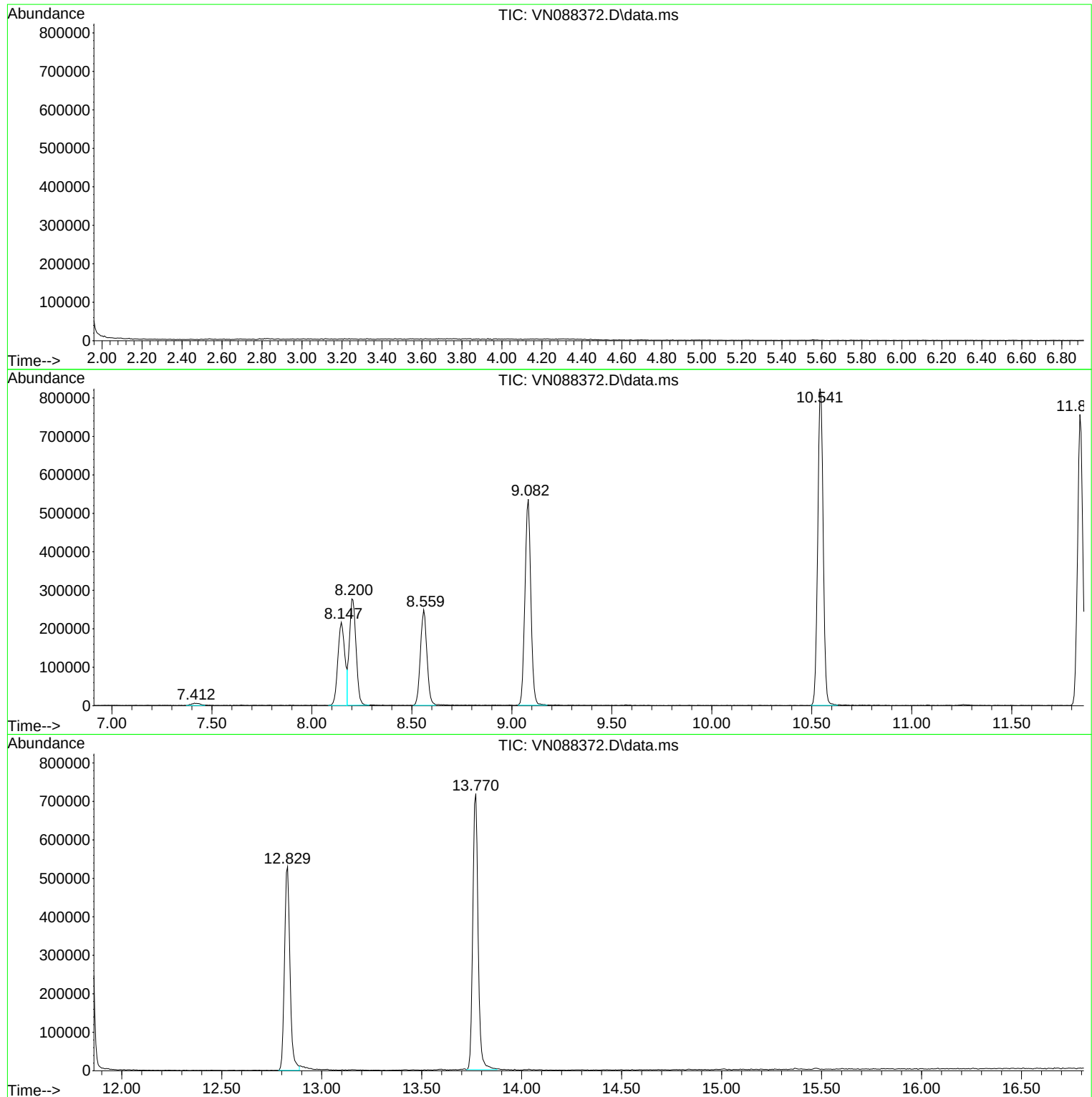
Sum of corrected areas: 8133289

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN120125\
Data File : VN088372.D
Acq On : 01 Dec 2025 10:53
Operator : JC\MD
Sample : VN1201WBL01
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 5 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
VN1201WBL01

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N112425W.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN120125\
Data File : VN088372.D
Acq On : 01 Dec 2025 10:53
Operator : JC\MD
Sample : VN1201WBL01
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 5 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
VN1201WBL01

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N112425W.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

No Library Search Compounds Detected

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN120125\
Data File : VN088372.D
Acq On : 01 Dec 2025 10:53
Operator : JC\MD
Sample : VN1201WBL01
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 5 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
VN1201WBL01

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N112425W.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
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Report of Analysis

Client: Remington & Vernick
 Project: Saddler Property
 Client Sample ID: VN1201WBS01
 Lab Sample ID: VN1201WBS01
 Analytical Method: 8260D
 Sample Wt/Vol: 5 mL

Level: LOW
 Final Vol: 5000 uL

Date Collected:
 Date Received:
 SDG No.: Q3743
 Matrix: Water
 % Solid: 0
 Test: VOC-TCLVOA-10

CAS Number	Parameter	Conc.	Qua.	DF	MDL	LOQ / CRQL	Units	Date Ana.	BatchID
TARGETS									
75-71-8	Dichlorodifluoromethane	18.2		1	0.22	1.00	ug/L	12/01/25 11:27	VN120125
74-87-3	Chloromethane	19.4		1	0.32	1.00	ug/L	12/01/25 11:27	VN120125
75-01-4	Vinyl Chloride	18.9		1	0.26	1.00	ug/L	12/01/25 11:27	VN120125
74-83-9	Bromomethane	19.7		1	1.40	5.00	ug/L	12/01/25 11:27	VN120125
75-00-3	Chloroethane	19.5		1	0.47	1.00	ug/L	12/01/25 11:27	VN120125
75-69-4	Trichlorofluoromethane	18.6		1	0.33	1.00	ug/L	12/01/25 11:27	VN120125
76-13-1	1,1,2-Trichlorotrifluoroethane	18.6		1	0.25	1.00	ug/L	12/01/25 11:27	VN120125
75-35-4	1,1-Dichloroethene	19.0		1	0.23	1.00	ug/L	12/01/25 11:27	VN120125
67-64-1	Acetone	87.9		1	1.50	5.00	ug/L	12/01/25 11:27	VN120125
75-15-0	Carbon Disulfide	19.4		1	0.21	1.00	ug/L	12/01/25 11:27	VN120125
1634-04-4	Methyl tert-butyl Ether	19.9		1	0.16	1.00	ug/L	12/01/25 11:27	VN120125
79-20-9	Methyl Acetate	19.1		1	0.27	1.00	ug/L	12/01/25 11:27	VN120125
75-09-2	Methylene Chloride	20.0		1	0.28	1.00	ug/L	12/01/25 11:27	VN120125
156-60-5	trans-1,2-Dichloroethene	19.6		1	0.23	1.00	ug/L	12/01/25 11:27	VN120125
75-34-3	1,1-Dichloroethane	19.8		1	0.23	1.00	ug/L	12/01/25 11:27	VN120125
110-82-7	Cyclohexane	17.9		1	1.50	5.00	ug/L	12/01/25 11:27	VN120125
78-93-3	2-Butanone	96.0		1	0.98	5.00	ug/L	12/01/25 11:27	VN120125
56-23-5	Carbon Tetrachloride	18.3		1	0.25	1.00	ug/L	12/01/25 11:27	VN120125
156-59-2	cis-1,2-Dichloroethene	20.1		1	0.19	1.00	ug/L	12/01/25 11:27	VN120125
74-97-5	Bromochloromethane	19.3		1	0.22	1.00	ug/L	12/01/25 11:27	VN120125
67-66-3	Chloroform	20.3		1	0.25	1.00	ug/L	12/01/25 11:27	VN120125
71-55-6	1,1,1-Trichloroethane	18.8		1	0.20	1.00	ug/L	12/01/25 11:27	VN120125
108-87-2	Methylcyclohexane	17.1		1	0.16	1.00	ug/L	12/01/25 11:27	VN120125
71-43-2	Benzene	19.4		1	0.15	1.00	ug/L	12/01/25 11:27	VN120125
107-06-2	1,2-Dichloroethane	19.6		1	0.22	1.00	ug/L	12/01/25 11:27	VN120125
79-01-6	Trichloroethene	18.8		1	0.090	1.00	ug/L	12/01/25 11:27	VN120125
78-87-5	1,2-Dichloropropane	19.8		1	0.20	1.00	ug/L	12/01/25 11:27	VN120125
75-27-4	Bromodichloromethane	19.5		1	0.22	1.00	ug/L	12/01/25 11:27	VN120125
108-10-1	4-Methyl-2-Pentanone	98.3		1	0.68	5.00	ug/L	12/01/25 11:27	VN120125
108-88-3	Toluene	19.5		1	0.14	1.00	ug/L	12/01/25 11:27	VN120125
10061-02-6	t-1,3-Dichloropropene	19.9		1	0.17	1.00	ug/L	12/01/25 11:27	VN120125
10061-01-5	cis-1,3-Dichloropropene	20.1		1	0.16	1.00	ug/L	12/01/25 11:27	VN120125
79-00-5	1,1,2-Trichloroethane	20.2		1	0.21	1.00	ug/L	12/01/25 11:27	VN120125
591-78-6	2-Hexanone	93.5		1	0.89	5.00	ug/L	12/01/25 11:27	VN120125
124-48-1	Dibromochloromethane	20.1		1	0.18	1.00	ug/L	12/01/25 11:27	VN120125
106-93-4	1,2-Dibromoethane	20.6		1	0.15	1.00	ug/L	12/01/25 11:27	VN120125
127-18-4	Tetrachloroethene	17.9		1	0.23	1.00	ug/L	12/01/25 11:27	VN120125
108-90-7	Chlorobenzene	19.5		1	0.12	1.00	ug/L	12/01/25 11:27	VN120125
100-41-4	Ethyl Benzene	18.6		1	0.13	1.00	ug/L	12/01/25 11:27	VN120125
179601-23-1	m/p-Xylenes	37.0		1	0.24	2.00	ug/L	12/01/25 11:27	VN120125
95-47-6	o-Xylene	18.3		1	0.12	1.00	ug/L	12/01/25 11:27	VN120125
100-42-5	Styrene	19.6		1	0.15	1.00	ug/L	12/01/25 11:27	VN120125
75-25-2	Bromoform	20.0		1	0.19	1.00	ug/L	12/01/25 11:27	VN120125



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

Report of Analysis

Client: Remington & Vernick
Project: Saddler Property
Client Sample ID: VN1201WBS01
Lab Sample ID: VN1201WBS01
Analytical Method: 8260D
Sample Wt/Vol: 5 mL

Level: LOW
Final Vol: 5000 uL

Date Collected:
Date Received:
SDG No.: Q3743
Matrix: Water
% Solid: 0
Test: VOC-TCLVOA-10

CAS Number	Parameter	Conc.	Qua.	DF	MDL	LOQ / CRQL	Units	Date Ana.	BatchID
98-82-8	Isopropylbenzene	18.6		1	0.12	1.00	ug/L	12/01/25 11:27	VN120125
79-34-5	1,1,2,2-Tetrachloroethane	20.1		1	0.26	1.00	ug/L	12/01/25 11:27	VN120125
541-73-1	1,3-Dichlorobenzene	19.0		1	0.16	1.00	ug/L	12/01/25 11:27	VN120125
106-46-7	1,4-Dichlorobenzene	18.5		1	0.19	1.00	ug/L	12/01/25 11:27	VN120125
95-50-1	1,2-Dichlorobenzene	19.4		1	0.16	1.00	ug/L	12/01/25 11:27	VN120125
96-12-8	1,2-Dibromo-3-Chloropropane	18.2		1	0.53	1.00	ug/L	12/01/25 11:27	VN120125
120-82-1	1,2,4-Trichlorobenzene	18.5		1	0.20	1.00	ug/L	12/01/25 11:27	VN120125
87-61-6	1,2,3-Trichlorobenzene	18.3		1	0.20	1.00	ug/L	12/01/25 11:27	VN120125
SURROGATES									
17060-07-0	1,2-Dichloroethane-d4	56.0			74 - 125	112%	SPK: 50		
1868-53-7	Dibromofluoromethane	55.1			75 - 124	110%	SPK: 50		
2037-26-5	Toluene-d8	54.9			86 - 113	110%	SPK: 50		
460-00-4	4-Bromofluorobenzene	53.3			77 - 121	107%	SPK: 50		
INTERNAL STANDARDS									
		Area Count							
363-72-4	Pentafluorobenzene	257000							
540-36-3	1,4-Difluorobenzene	487000							
3114-55-4	Chlorobenzene-d5	435000							
3855-82-1	1,4-Dichlorobenzene-d4	207000							

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\VN120125\
 Data File : VN088373.D
 Acq On : 01 Dec 2025 11:27
 Operator : JC\MD
 Sample : VN1201WBS01
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 VN1201WBS01

Manual Integrations
 APPROVED

Reviewed By : John Carlone 12/02/2025
 Supervised By : Mahesh Dadoda 12/02/2025

Quant Time: Dec 02 00:10:02 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	256678	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.083	114	487472	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.841	117	435019	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.770	152	207303	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.559	65	271013	56.049	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery	=	112.100%	
35) Dibromofluoromethane	8.147	113	186190	55.116	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery	=	110.240%	
50) Toluene-d8	10.541	98	690335	54.922	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery	=	109.840%	
62) 4-Bromofluorobenzene	12.829	95	230018	53.334	ug/l	0.00
Spiked Amount 50.000	Range 77 - 121		Recovery	=	106.660%	
Target Compounds						Qvalue
2) Dichlorodifluoromethane	2.142	85	70210	18.211	ug/l	99
3) Chloromethane	2.383	50	80214	19.372	ug/l	94
4) Vinyl Chloride	2.536	62	79487	18.870	ug/l	99
5) Bromomethane	2.989	94	39940	19.700	ug/l	96
6) Chloroethane	3.142	64	50911	19.530	ug/l	99
7) Trichlorofluoromethane	3.512	101	109953	18.572	ug/l	94
8) Diethyl Ether	3.965	74	43004	19.352	ug/l	94
9) 1,1,2-Trichlorotrifluo...	4.371	101	58508	18.563	ug/l	97
10) Methyl Iodide	4.583	142	59901	19.198	ug/l	98
11) Tert butyl alcohol	5.512	59	67525	87.469	ug/l	97
12) 1,1-Dichloroethene	4.336	96	59547	19.015	ug/l	87
13) Acrolein	4.177	56	88765	120.102	ug/l	97
14) Allyl chloride	5.018	41	108153	18.633	ug/l	98
15) Acrylonitrile	5.706	53	218028	97.888	ug/l	98
16) Acetone	4.418	43	153779	87.939	ug/l	94
17) Carbon Disulfide	4.706	76	198994	19.446	ug/l	100
18) Methyl Acetate	5.018	43	99423	19.100	ug/l	98
19) Methyl tert-butyl Ether	5.783	73	233707	19.919	ug/l	100
20) Methylene Chloride	5.271	84	74861	20.009	ug/l	83
21) trans-1,2-Dichloroethene	5.771	96	68123	19.638	ug/l	89
22) Diisopropyl ether	6.653	45	255505	20.619	ug/l	93
23) Vinyl Acetate	6.589	43	1003026m	100.465	ug/l	
24) 1,1-Dichloroethane	6.547	63	139057	19.816	ug/l	100
25) 2-Butanone	7.465	43	269940	96.022	ug/l	97
26) 2,2-Dichloropropane	7.471	77	115800	19.118	ug/l	98
27) cis-1,2-Dichloroethene	7.465	96	80954	20.080	ug/l	98
28) Bromochloromethane	7.789	49	67630	19.346	ug/l #	98
29) Tetrahydrofuran	7.818	42	200576	99.471	ug/l	99
30) Chloroform	7.941	83	140173	20.329	ug/l	100
31) Cyclohexane	8.241	56	114116	17.945	ug/l	98
32) 1,1,1-Trichloroethane	8.147	97	114251	18.845	ug/l	96
36) 1,1-Dichloropropene	8.353	75	93124	17.862	ug/l	99
37) Ethyl Acetate	7.547	43	129007	18.819	ug/l	100
38) Carbon Tetrachloride	8.341	117	94912	18.287	ug/l	96
39) Methylcyclohexane	9.583	83	95473	17.120	ug/l	97
40) Benzene	8.588	78	296136	19.353	ug/l	99

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN120125\
 Data File : VN088373.D
 Acq On : 01 Dec 2025 11:27
 Operator : JC\MD
 Sample : VN1201WBS01
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 VN1201WBS01

Manual Integrations
 APPROVED

Reviewed By : John Carlone 12/02/2025
 Supervised By : Mahesh Dadoda 12/02/2025

Quant Time: Dec 02 00:10:02 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	63234	18.624	ug/l	98
42) 1,2-Dichloroethane	8.653	62	114468	19.607	ug/l	98
43) Isopropyl Acetate	8.671	43	195071	18.624	ug/l	98
44) Trichloroethene	9.330	130	67826	18.759	ug/l	95
45) 1,2-Dichloropropane	9.600	63	77770	19.837	ug/l	93
46) Dibromomethane	9.688	93	55251	19.747	ug/l	98
47) Bromodichloromethane	9.871	83	111551	19.540	ug/l	92
48) Methyl methacrylate	9.659	41	87729	18.357	ug/l	97
49) 1,4-Dioxane	9.683	88	21595	355.324	ug/l #	99
51) 4-Methyl-2-Pentanone	10.424	43	606997	98.320	ug/l	100
52) Toluene	10.606	92	172323	19.503	ug/l	99
53) t-1,3-Dichloropropene	10.812	75	116655	19.908	ug/l	98
54) cis-1,3-Dichloropropene	10.294	75	123675	20.070	ug/l	99
55) 1,1,2-Trichloroethane	10.994	97	68939	20.166	ug/l	96
56) Ethyl methacrylate	10.859	69	106640	19.749	ug/l	96
57) 1,3-Dichloropropane	11.141	76	122169	19.810	ug/l	98
58) 2-Chloroethyl Vinyl ether	10.135	63	278967	89.610	ug/l	97
59) 2-Hexanone	11.177	43	363554	93.479	ug/l	98
60) Dibromochloromethane	11.335	129	78151	20.054	ug/l	97
61) 1,2-Dibromoethane	11.447	107	71254	20.583	ug/l	98
64) Tetrachloroethene	11.082	164	61757	17.907	ug/l	96
65) Chlorobenzene	11.871	112	190129	19.458	ug/l	97
66) 1,1,1,2-Tetrachloroethane	11.935	131	65977	20.020	ug/l	98
67) Ethyl Benzene	11.941	91	316848	18.553	ug/l	99
68) m/p-Xylenes	12.047	106	234320	37.005	ug/l	97
69) o-Xylene	12.376	106	110515	18.307	ug/l	100
70) Styrene	12.388	104	191890	19.577	ug/l	99
71) Bromoform	12.559	173	48189	19.963	ug/l #	99
73) Isopropylbenzene	12.676	105	285585	18.632	ug/l	99
74) N-amyl acetate	12.682	43	103450m	19.818	ug/l	
75) 1,1,2,2-Tetrachloroethane	12.912	83	99427	20.065	ug/l	97
76) 1,2,3-Trichloropropane	12.971	75	95762m	19.723	ug/l	
77) Bromobenzene	12.959	156	70858	20.205	ug/l	97
78) n-propylbenzene	13.012	91	349255	18.625	ug/l	100
79) 2-Chlorotoluene	13.100	91	213346	18.631	ug/l	98
80) 1,3,5-Trimethylbenzene	13.147	105	239698	18.872	ug/l	98
81) trans-1,4-Dichloro-2-b...	12.718	75	31603	18.465	ug/l #	98
82) 4-Chlorotoluene	13.200	91	222136	19.531	ug/l	99
83) tert-Butylbenzene	13.412	119	195850	18.139	ug/l	99
84) 1,2,4-Trimethylbenzene	13.459	105	244186	19.342	ug/l	99
85) sec-Butylbenzene	13.594	105	282563	18.217	ug/l	100
86) p-Isopropyltoluene	13.706	119	237355	18.661	ug/l	99
87) 1,3-Dichlorobenzene	13.712	146	129958	19.036	ug/l	99
88) 1,4-Dichlorobenzene	13.788	146	131944	18.451	ug/l	97
89) n-Butylbenzene	14.029	91	220480	17.853	ug/l	100
90) Hexachloroethane	14.306	117	42679	18.368	ug/l	95
91) 1,2-Dichlorobenzene	14.082	146	124887	19.439	ug/l	98
92) 1,2-Dibromo-3-Chloropr...	14.694	75	21886	18.204	ug/l	99
93) 1,2,4-Trichlorobenzene	15.365	180	70611	18.498	ug/l	98
94) Hexachlorobutadiene	15.470	225	24611	17.942	ug/l	99
95) Naphthalene	15.612	128	234599	17.736	ug/l	99
96) 1,2,3-Trichlorobenzene	15.812	180	68004	18.306	ug/l	99

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN120125\
Data File : VN088373.D
Acq On : 01 Dec 2025 11:27
Operator : JC\MD
Sample : VN1201WBS01
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
VN1201WBS01

Manual Integrations
APPROVED

Reviewed By :John Carlone 12/02/2025
Supervised By :Mahesh Dadoda 12/02/2025

Quant Time: Dec 02 00:10:02 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)
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(#) = qualifier out of range (m) = manual integration (+) = signals summed

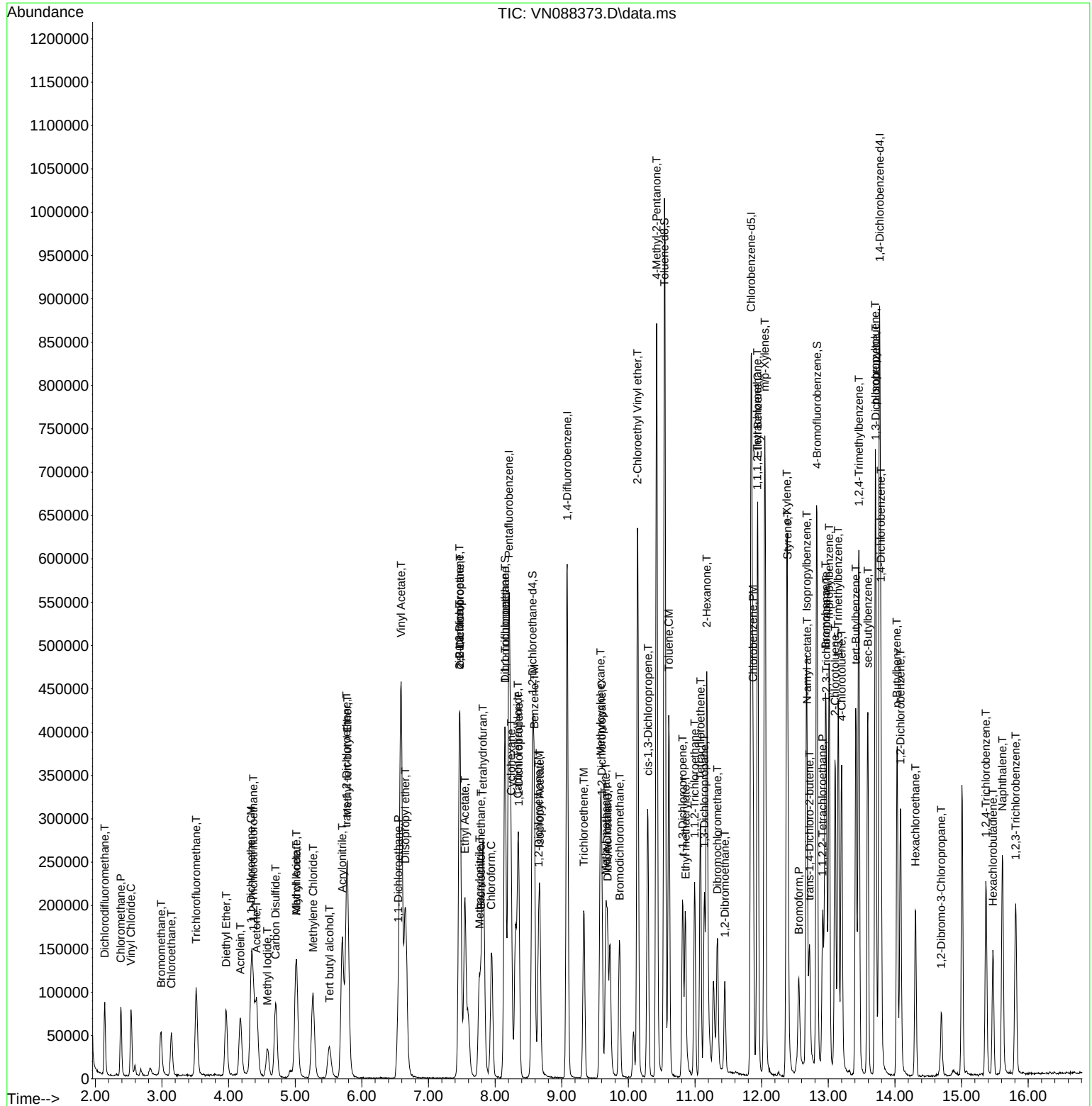
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Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN120125\  
Data File : VN088373.D  
Acq On    : 01 Dec 2025  11:27  
Operator  : JC\MD  
Sample    : VN1201WBS01  
Misc      : 5.0mL/MSVOA_N/WATER  
ALS Vial  : 6    Sample Multiplier: 1
```

Instrument :
MSVOA_N
ClientSampleId :
VN1201WBS01

Manual Integrations

Reviewed By :John Carlone 12/02/2025
Supervised By :Mahesh Dadoda 12/02/2025

Quant Time: Dec 02 00:10:02 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



Report of Analysis

Client: Remington & Vernick
Project: Saddler Property
Client Sample ID: VN1201WBSD01
Lab Sample ID: VN1201WBSD01
Analytical Method: 8260D
Sample Wt/Vol: 5 mL

Level: LOW
Final Vol: 5000 uL

Date Collected:
Date Received:
SDG No.: Q3743
Matrix: Water
% Solid: 0
Test: VOC-TCLVOA-10

CAS Number	Parameter	Conc.	Qua.	DF	MDL	LOQ / CRQL	Units	Date Ana.	BatchID
TARGETS									
75-71-8	Dichlorodifluoromethane	18.9		1	0.22	1.00	ug/L	12/01/25 12:10	VN120125
74-87-3	Chloromethane	19.4		1	0.32	1.00	ug/L	12/01/25 12:10	VN120125
75-01-4	Vinyl Chloride	18.7		1	0.26	1.00	ug/L	12/01/25 12:10	VN120125
74-83-9	Bromomethane	20.2		1	1.40	5.00	ug/L	12/01/25 12:10	VN120125
75-00-3	Chloroethane	19.3		1	0.47	1.00	ug/L	12/01/25 12:10	VN120125
75-69-4	Trichlorofluoromethane	18.7		1	0.33	1.00	ug/L	12/01/25 12:10	VN120125
76-13-1	1,1,2-Trichlorotrifluoroethane	18.5		1	0.25	1.00	ug/L	12/01/25 12:10	VN120125
75-35-4	1,1-Dichloroethene	18.8		1	0.23	1.00	ug/L	12/01/25 12:10	VN120125
67-64-1	Acetone	87.0		1	1.50	5.00	ug/L	12/01/25 12:10	VN120125
75-15-0	Carbon Disulfide	18.8		1	0.21	1.00	ug/L	12/01/25 12:10	VN120125
1634-04-4	Methyl tert-butyl Ether	19.4		1	0.16	1.00	ug/L	12/01/25 12:10	VN120125
79-20-9	Methyl Acetate	19.0		1	0.27	1.00	ug/L	12/01/25 12:10	VN120125
75-09-2	Methylene Chloride	20.2		1	0.28	1.00	ug/L	12/01/25 12:10	VN120125
156-60-5	trans-1,2-Dichloroethene	19.6		1	0.23	1.00	ug/L	12/01/25 12:10	VN120125
75-34-3	1,1-Dichloroethane	19.1		1	0.23	1.00	ug/L	12/01/25 12:10	VN120125
110-82-7	Cyclohexane	18.5		1	1.50	5.00	ug/L	12/01/25 12:10	VN120125
78-93-3	2-Butanone	96.2		1	0.98	5.00	ug/L	12/01/25 12:10	VN120125
56-23-5	Carbon Tetrachloride	18.8		1	0.25	1.00	ug/L	12/01/25 12:10	VN120125
156-59-2	cis-1,2-Dichloroethene	19.4		1	0.19	1.00	ug/L	12/01/25 12:10	VN120125
74-97-5	Bromochloromethane	18.5		1	0.22	1.00	ug/L	12/01/25 12:10	VN120125
67-66-3	Chloroform	19.6		1	0.25	1.00	ug/L	12/01/25 12:10	VN120125
71-55-6	1,1,1-Trichloroethane	19.0		1	0.20	1.00	ug/L	12/01/25 12:10	VN120125
108-87-2	Methylcyclohexane	18.5		1	0.16	1.00	ug/L	12/01/25 12:10	VN120125
71-43-2	Benzene	19.1		1	0.15	1.00	ug/L	12/01/25 12:10	VN120125
107-06-2	1,2-Dichloroethane	19.4		1	0.22	1.00	ug/L	12/01/25 12:10	VN120125
79-01-6	Trichloroethene	19.1		1	0.090	1.00	ug/L	12/01/25 12:10	VN120125
78-87-5	1,2-Dichloropropane	19.2		1	0.20	1.00	ug/L	12/01/25 12:10	VN120125
75-27-4	Bromodichloromethane	19.3		1	0.22	1.00	ug/L	12/01/25 12:10	VN120125
108-10-1	4-Methyl-2-Pentanone	98.5		1	0.68	5.00	ug/L	12/01/25 12:10	VN120125
108-88-3	Toluene	19.8		1	0.14	1.00	ug/L	12/01/25 12:10	VN120125
10061-02-6	t-1,3-Dichloropropene	19.1		1	0.17	1.00	ug/L	12/01/25 12:10	VN120125
10061-01-5	cis-1,3-Dichloropropene	19.5		1	0.16	1.00	ug/L	12/01/25 12:10	VN120125
79-00-5	1,1,2-Trichloroethane	20.0		1	0.21	1.00	ug/L	12/01/25 12:10	VN120125
591-78-6	2-Hexanone	94.4		1	0.89	5.00	ug/L	12/01/25 12:10	VN120125
124-48-1	Dibromochloromethane	20.3		1	0.18	1.00	ug/L	12/01/25 12:10	VN120125
106-93-4	1,2-Dibromoethane	19.9		1	0.15	1.00	ug/L	12/01/25 12:10	VN120125
127-18-4	Tetrachloroethene	19.0		1	0.23	1.00	ug/L	12/01/25 12:10	VN120125
108-90-7	Chlorobenzene	19.8		1	0.12	1.00	ug/L	12/01/25 12:10	VN120125
100-41-4	Ethyl Benzene	18.9		1	0.13	1.00	ug/L	12/01/25 12:10	VN120125
179601-23-1	m/p-Xylenes	37.8		1	0.24	2.00	ug/L	12/01/25 12:10	VN120125
95-47-6	o-Xylene	19.2		1	0.12	1.00	ug/L	12/01/25 12:10	VN120125
100-42-5	Styrene	19.4		1	0.15	1.00	ug/L	12/01/25 12:10	VN120125
75-25-2	Bromoform	20.1		1	0.19	1.00	ug/L	12/01/25 12:10	VN120125



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

Report of Analysis

Client: Remington & Vernick
Project: Saddler Property
Client Sample ID: VN1201WBSD01
Lab Sample ID: VN1201WBSD01
Analytical Method: 8260D
Sample Wt/Vol: 5 mL

Level: LOW
Final Vol: 5000 uL

Date Collected:
Date Received:
SDG No.: Q3743
Matrix: Water
% Solid: 0
Test: VOC-TCLVOA-10

CAS Number	Parameter	Conc.	Qua.	DF	MDL	LOQ / CRQL	Units	Date Ana.	BatchID
98-82-8	Isopropylbenzene	19.3		1	0.12	1.00	ug/L	12/01/25 12:10	VN120125
79-34-5	1,1,2,2-Tetrachloroethane	19.5		1	0.26	1.00	ug/L	12/01/25 12:10	VN120125
541-73-1	1,3-Dichlorobenzene	19.5		1	0.16	1.00	ug/L	12/01/25 12:10	VN120125
106-46-7	1,4-Dichlorobenzene	19.0		1	0.19	1.00	ug/L	12/01/25 12:10	VN120125
95-50-1	1,2-Dichlorobenzene	19.0		1	0.16	1.00	ug/L	12/01/25 12:10	VN120125
96-12-8	1,2-Dibromo-3-Chloropropane	18.5		1	0.53	1.00	ug/L	12/01/25 12:10	VN120125
120-82-1	1,2,4-Trichlorobenzene	18.0		1	0.20	1.00	ug/L	12/01/25 12:10	VN120125
87-61-6	1,2,3-Trichlorobenzene	19.1		1	0.20	1.00	ug/L	12/01/25 12:10	VN120125

SURROGATES

17060-07-0	1,2-Dichloroethane-d4	52.2			74 - 125	104%	SPK: 50
1868-53-7	Dibromofluoromethane	52.3			75 - 124	105%	SPK: 50
2037-26-5	Toluene-d8	53.2			86 - 113	106%	SPK: 50
460-00-4	4-Bromofluorobenzene	51.1			77 - 121	102%	SPK: 50

INTERNAL STANDARDS

Area Count

363-72-4	Pentafluorobenzene	257000
540-36-3	1,4-Difluorobenzene	473000
3114-55-4	Chlorobenzene-d5	413000
3855-82-1	1,4-Dichlorobenzene-d4	197000

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\VN120125\
 Data File : VN088374.D
 Acq On : 01 Dec 2025 12:10
 Operator : JC\MD
 Sample : VN1201WBSD01
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 VN1201WBSD01

Manual Integrations
 APPROVED

Reviewed By : John Carlone 12/02/2025
 Supervised By : Mahesh Dadoda 12/02/2025

Quant Time: Dec 02 00:11:04 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	257123	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.082	114	472668	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.841	117	413438	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.770	152	196888	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.559	65	252803	52.192	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery = 104.380%			
35) Dibromofluoromethane	8.147	113	171251	52.281	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery = 104.560%			
50) Toluene-d8	10.541	98	648629	53.220	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery = 106.440%			
62) 4-Bromofluorobenzene	12.829	95	213607	51.080	ug/l	0.00
Spiked Amount 50.000	Range 77 - 121		Recovery = 102.160%			
Target Compounds					Qvalue	
2) Dichlorodifluoromethane	2.142	85	72877	18.870	ug/l	98
3) Chloromethane	2.383	50	80587	19.429	ug/l	94
4) Vinyl Chloride	2.536	62	79060	18.736	ug/l	99
5) Bromomethane	2.989	94	40980	20.178	ug/l	99
6) Chloroethane	3.148	64	50269	19.250	ug/l	90
7) Trichlorofluoromethane	3.512	101	110683	18.663	ug/l	99
8) Diethyl Ether	3.965	74	42809	19.231	ug/l	98
9) 1,1,2-Trichlorotrifluo...	4.377	101	58334	18.476	ug/l	98
10) Methyl Iodide	4.583	142	60569	19.379	ug/l	96
11) Tert butyl alcohol	5.512	59	68039	87.983	ug/l #	81
12) 1,1-Dichloroethene	4.347	96	58831	18.754	ug/l	84
13) Acrolein	4.177	56	87619	118.346	ug/l	97
14) Allyl chloride	5.012	41	104505	17.974	ug/l	98
15) Acrylonitrile	5.706	53	213646	95.755	ug/l	98
16) Acetone	4.418	43	152332	86.961	ug/l	98
17) Carbon Disulfide	4.706	76	192813	18.810	ug/l	100
18) Methyl Acetate	5.018	43	99171	19.019	ug/l	99
19) Methyl tert-butyl Ether	5.783	73	228498	19.442	ug/l	95
20) Methylene Chloride	5.259	84	75693	20.196	ug/l	83
21) trans-1,2-Dichloroethene	5.771	96	68139	19.608	ug/l	92
22) Diisopropyl ether	6.653	45	244953	19.733	ug/l	96
23) Vinyl Acetate	6.588	43	972820m	97.271	ug/l	
24) 1,1-Dichloroethane	6.553	63	134006	19.063	ug/l	95
25) 2-Butanone	7.465	43	270811	96.165	ug/l	97
26) 2,2-Dichloropropane	7.477	77	112288	18.506	ug/l	95
27) cis-1,2-Dichloroethene	7.465	96	78314	19.392	ug/l	96
28) Bromochloromethane	7.794	49	64631	18.456	ug/l	99
29) Tetrahydrofuran	7.824	42	193749	95.919	ug/l	99
30) Chloroform	7.947	83	135282	19.585	ug/l	96
31) Cyclohexane	8.235	56	117687	18.475	ug/l	98
32) 1,1,1-Trichloroethane	8.147	97	115347	18.992	ug/l	95
36) 1,1-Dichloropropene	8.353	75	89956	17.794	ug/l	98
37) Ethyl Acetate	7.547	43	123427	18.569	ug/l	98
38) Carbon Tetrachloride	8.347	117	94617	18.801	ug/l	96
39) Methylcyclohexane	9.576	83	99985	18.491	ug/l	93
40) Benzene	8.588	78	284116	19.149	ug/l	98

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN120125\
 Data File : VN088374.D
 Acq On : 01 Dec 2025 12:10
 Operator : JC\MD
 Sample : VN1201WBSD01
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 VN1201WBSD01

Manual Integrations
 APPROVED

Reviewed By : John Carlone 12/02/2025
 Supervised By : Mahesh Dadoda 12/02/2025

Quant Time: Dec 02 00:11:04 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	63105	19.168	ug/l	98
42) 1,2-Dichloroethane	8.647	62	109703	19.380	ug/l	98
43) Isopropyl Acetate	8.671	43	192472	18.952	ug/l	99
44) Trichloroethene	9.329	130	66801	19.054	ug/l	90
45) 1,2-Dichloropropane	9.600	63	72843	19.162	ug/l	100
46) Dibromomethane	9.688	93	54401	20.053	ug/l	98
47) Bromodichloromethane	9.865	83	106876	19.308	ug/l #	93
48) Methyl methacrylate	9.659	41	84615	18.260	ug/l	99
49) 1,4-Dioxane	9.676	88	22017	373.614	ug/l #	99
51) 4-Methyl-2-Pentanone	10.424	43	589357	98.453	ug/l	99
52) Toluene	10.606	92	170024	19.846	ug/l	98
53) t-1,3-Dichloropropene	10.818	75	108520	19.099	ug/l	99
54) cis-1,3-Dichloropropene	10.288	75	116758	19.541	ug/l	97
55) 1,1,2-Trichloroethane	10.994	97	66233	19.982	ug/l	95
56) Ethyl methacrylate	10.853	69	104863	20.028	ug/l	97
57) 1,3-Dichloropropane	11.141	76	120668	20.179	ug/l	99
58) 2-Chloroethyl Vinyl ether	10.141	63	276569	91.622	ug/l	99
59) 2-Hexanone	11.176	43	355826	94.358	ug/l	98
60) Dibromochloromethane	11.335	129	76699	20.298	ug/l	98
61) 1,2-Dibromoethane	11.447	107	66826	19.909	ug/l	97
64) Tetrachloroethene	11.076	164	62208	18.980	ug/l	95
65) Chlorobenzene	11.870	112	183680	19.779	ug/l	96
66) 1,1,1,2-Tetrachloroethane	11.941	131	61851	19.747	ug/l	99
67) Ethyl Benzene	11.941	91	306102	18.859	ug/l	97
68) m/p-Xylenes	12.053	106	227659	37.830	ug/l	99
69) o-Xylene	12.376	106	109896	19.155	ug/l	95
70) Styrene	12.388	104	181083	19.438	ug/l	99
71) Bromoform	12.559	173	46125	20.106	ug/l #	99
73) Isopropylbenzene	12.670	105	280946	19.299	ug/l	100
74) N-amyl acetate	12.682	43	97186m	19.603	ug/l	
75) 1,1,2,2-Tetrachloroethane	12.917	83	91931	19.534	ug/l	97
76) 1,2,3-Trichloropropane	12.970	75	93113m	20.192	ug/l	
77) Bromobenzene	12.959	156	65049	19.529	ug/l	96
78) n-propylbenzene	13.012	91	348515	19.569	ug/l	100
79) 2-Chlorotoluene	13.100	91	204279	18.783	ug/l	99
80) 1,3,5-Trimethylbenzene	13.153	105	234071	19.404	ug/l	99
81) trans-1,4-Dichloro-2-b...	12.717	75	28199	17.348	ug/l #	96
82) 4-Chlorotoluene	13.200	91	211139	19.546	ug/l	97
83) tert-Butylbenzene	13.412	119	192736	18.795	ug/l	99
84) 1,2,4-Trimethylbenzene	13.459	105	235984	19.682	ug/l	100
85) sec-Butylbenzene	13.594	105	283289	19.230	ug/l	100
86) p-Isopropyltoluene	13.706	119	228290	18.898	ug/l	98
87) 1,3-Dichlorobenzene	13.712	146	126657	19.534	ug/l	98
88) 1,4-Dichlorobenzene	13.788	146	129333	19.042	ug/l	98
89) n-Butylbenzene	14.029	91	219415	18.707	ug/l	99
90) Hexachloroethane	14.306	117	41810	18.946	ug/l	96
91) 1,2-Dichlorobenzene	14.082	146	115702	18.962	ug/l	99
92) 1,2-Dibromo-3-Chloropr...	14.700	75	21103	18.481	ug/l	98
93) 1,2,4-Trichlorobenzene	15.370	180	65316	18.016	ug/l	95
94) Hexachlorobutadiene	15.470	225	24556	18.849	ug/l	94
95) Naphthalene	15.611	128	231154	18.400	ug/l	99
96) 1,2,3-Trichlorobenzene	15.811	180	67438	19.114	ug/l	98

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN120125\
Data File : VN088374.D
Acq On : 01 Dec 2025 12:10
Operator : JC\MD
Sample : VN1201WBSD01
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
VN1201WBSD01

Manual Integrations
APPROVED

Reviewed By :John Carlone 12/02/2025
Supervised By :Mahesh Dadoda 12/02/2025

Quant Time: Dec 02 00:11:04 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)
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(#) = qualifier out of range (m) = manual integration (+) = signals summed

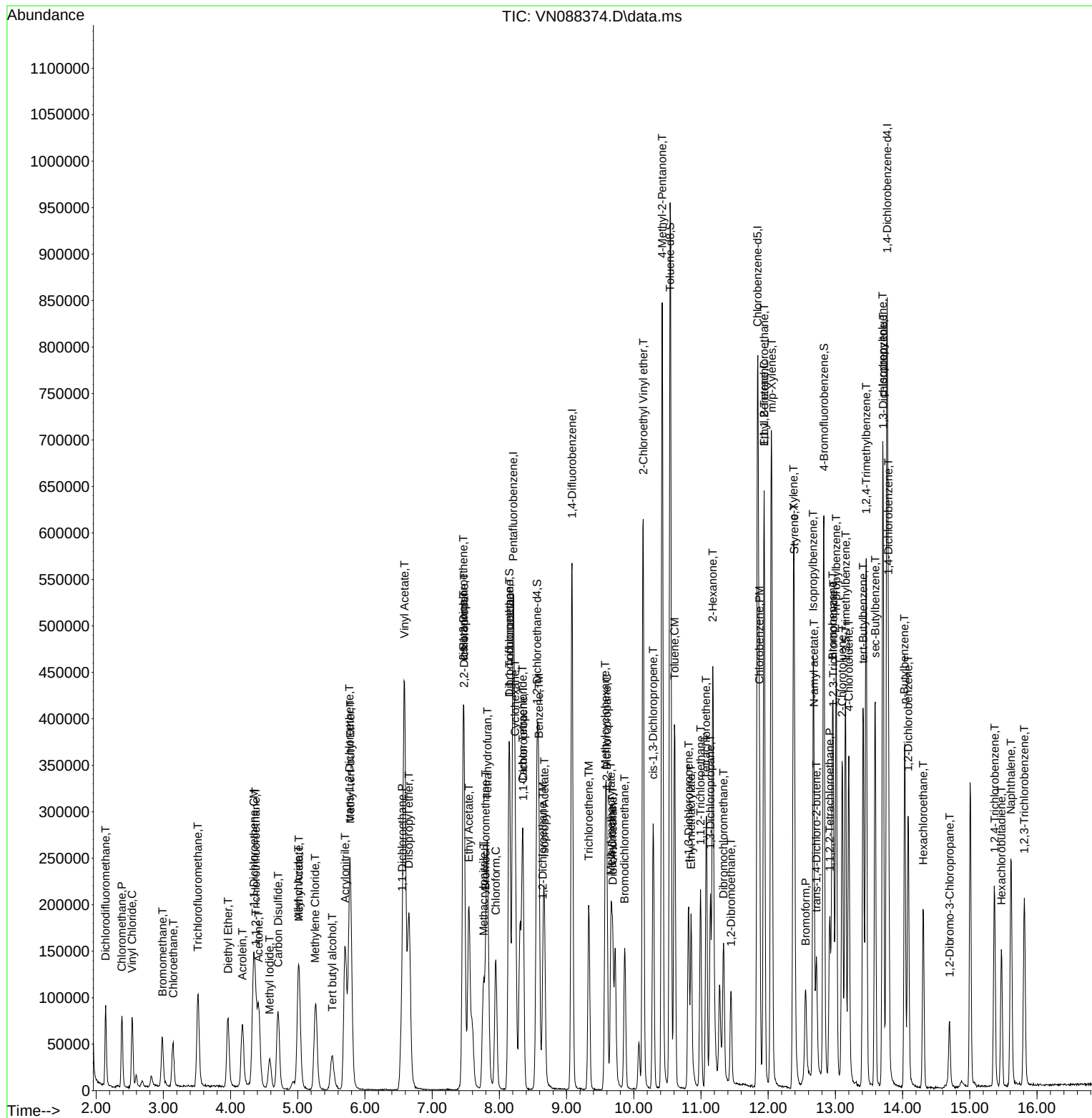
```
Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN120125\
Data File : VN088374.D
Acq On    : 01 Dec 2025  12:10
Operator  : JC\MD
Sample    : VN1201WBSD01
Misc      : 5.0mL/MSVOA_N/WATER
ALS Vial  : 7      Sample Multiplier: 1
```

Instrument :
MSVOA_N
ClientSampleId :
VN1201WBSD01

Manual Integrations APPROVED

Reviewed By :John Carlone 12/02/2025
Supervised By :Mahesh Dadoda 12/02/2025

Quant Time: Dec 02 00:11:04 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
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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:	VN120125	Instrument	MSVOA_n
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Sample ID	File ID	Parameter	Review By	Review On	Supervised By	Supervised On	Reason
VSTDCCC050	VN088370.D	1,2,3-Trichloropropane	JOHN	12/2/2025 7:49:00 AM	MMDadoda	12/2/2025 1:11:28 PM	Peak Integrated by Software
VSTDCCC050	VN088370.D	Methacrylonitrile	JOHN	12/2/2025 7:49:00 AM	MMDadoda	12/2/2025 1:11:28 PM	Peak Integrated by Software
VSTDCCC050	VN088370.D	N-amyl acetate	JOHN	12/2/2025 7:49:00 AM	MMDadoda	12/2/2025 1:11:28 PM	Peak Integrated by Software
VSTDCCC050	VN088370.D	Vinyl Acetate	JOHN	12/2/2025 7:49:00 AM	MMDadoda	12/2/2025 1:11:28 PM	Peak Integrated by Software
VN1201WBS01	VN088373.D	1,2,3-Trichloropropane	JOHN	12/2/2025 7:49:05 AM	MMDadoda	12/2/2025 1:11:32 PM	Peak Integrated by Software
VN1201WBS01	VN088373.D	N-amyl acetate	JOHN	12/2/2025 7:49:05 AM	MMDadoda	12/2/2025 1:11:32 PM	Peak Integrated by Software
VN1201WBS01	VN088373.D	Vinyl Acetate	JOHN	12/2/2025 7:49:05 AM	MMDadoda	12/2/2025 1:11:32 PM	Peak Integrated by Software
VN1201WBSD01	VN088374.D	1,2,3-Trichloropropane	JOHN	12/2/2025 7:49:10 AM	MMDadoda	12/2/2025 1:11:38 PM	Peak Integrated by Software
VN1201WBSD01	VN088374.D	N-amyl acetate	JOHN	12/2/2025 7:49:10 AM	MMDadoda	12/2/2025 1:11:38 PM	Peak Integrated by Software
VN1201WBSD01	VN088374.D	Vinyl Acetate	JOHN	12/2/2025 7:49:10 AM	MMDadoda	12/2/2025 1:11:38 PM	Peak Integrated by Software
VSTDCCC050	VN088394.D	1,2,3-Trichloropropane	JOHN	12/2/2025 7:49:42 AM	MMDadoda	12/2/2025 1:12:07 PM	Peak Integrated by Software
VSTDCCC050	VN088394.D	N-amyl acetate	JOHN	12/2/2025 7:49:42 AM	MMDadoda	12/2/2025 1:12:07 PM	Peak Integrated by Software
VSTDCCC050	VN088394.D	Vinyl Acetate	JOHN	12/2/2025 7:49:42 AM	MMDadoda	12/2/2025 1:12:07 PM	Peak Integrated by Software



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

Manual Integration Report

Sequence:

VN120125

Instrument

MSVOA_n

Sample ID

File ID

Parameter

Review By

Review On

Supervised
By

Supervised On

Reason

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 # VN120125

Review By	Semsettin Yesilyurt	Review On	12/4/2025 10:27:11 AM		
Supervise By	John Carlone	Supervise On	12/4/2025 10:36:07 AM		
SubDirectory	VN120125	HP Acquire Method	MSVOA_N	HP Processing Method	82N112425W.M
STD. NAME		STD REF.#			
Tune/Reschk Initial Calibration Stds		VP136451			
CCC Internal Standard/PEM ICV/I.BLK Surrogate Standard MS/MSD Standard LCS Standard		VP136452,VP136453			

Sr#	SampleId	Data File Name	Date-Time	Operator	Status
1	BFB	VN088369.D	01 Dec 2025 08:49	JC\MD	Ok
2	VSTDCCC050	VN088370.D	01 Dec 2025 09:58	JC\MD	Ok,M
3	VN1201MBL01	VN088371.D	01 Dec 2025 10:33	JC\MD	Ok
4	VN1201WBL01	VN088372.D	01 Dec 2025 10:53	JC\MD	Ok
5	VN1201WBS01	VN088373.D	01 Dec 2025 11:27	JC\MD	Ok,M
6	VN1201WBSD01	VN088374.D	01 Dec 2025 12:10	JC\MD	Ok,M
7	Q3715-02	VN088375.D	01 Dec 2025 12:30	JC\MD	Ok,M
8	Q3741-07	VN088376.D	01 Dec 2025 12:51	JC\MD	Ok
9	Q3716-01	VN088377.D	01 Dec 2025 13:11	JC\MD	Dilution
10	Q3716-03	VN088378.D	01 Dec 2025 13:32	JC\MD	Not Ok
11	Q3716-04	VN088379.D	01 Dec 2025 13:52	JC\MD	Dilution
12	Q3715-01	VN088380.D	01 Dec 2025 14:13	JC\MD	ReRun
13	Q3716-02	VN088381.D	01 Dec 2025 14:33	JC\MD	Ok
14	Q3738-01	VN088382.D	01 Dec 2025 14:54	JC\MD	Not Ok
15	Q3738-03	VN088383.D	01 Dec 2025 15:15	JC\MD	Ok
16	Q3738-02	VN088384.D	01 Dec 2025 15:35	JC\MD	Ok,M
17	Q3745-01	VN088385.D	01 Dec 2025 15:56	JC\MD	ReRun
18	Q3745-02	VN088386.D	01 Dec 2025 16:16	JC\MD	Ok
19	Q3745-03	VN088387.D	01 Dec 2025 16:37	JC\MD	Ok
20	Q3745-04	VN088388.D	01 Dec 2025 16:58	JC\MD	Ok
21	Q3743-03	VN088389.D	01 Dec 2025 17:18	JC\MD	Ok

Instrument ID: **MSVOA_N**

Daily Analysis Runlog For Sequence/QCBatch ID # VN120125

Review By	Semsettin Yesilyurt	Review On	12/4/2025 10:27:11 AM		
Supervise By	John Carlone	Supervise On	12/4/2025 10:36:07 AM		
SubDirectory	VN120125	HP Acquire Method	MSVOA_N	HP Processing Method	82N112425W.M
STD. NAME		STD REF.#			
Tune/Reschk Initial Calibration Stds		VP136451			
CCC Internal Standard/PEM ICV/I.BLK Surrogate Standard MS/MSD Standard LCS Standard		VP136452,VP136453			

22	Q3743-04	VN088390.D	01 Dec 2025 17:39	JC\MD	Ok
23	Q3743-05	VN088391.D	01 Dec 2025 18:00	JC\MD	Not Ok
24	Q3743-01	VN088392.D	01 Dec 2025 18:20	JC\MD	Ok
25	Q3743-02	VN088393.D	01 Dec 2025 18:41	JC\MD	Ok
26	VSTDCCC050	VN088394.D	01 Dec 2025 19:01	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 # VN120125

Review By	Semsettin Yesilyurt	Review On	12/4/2025 10:27:11 AM
Supervise By	John Carlone	Supervise On	12/4/2025 10:36:07 AM
SubDirectory	VN120125	HP Acquire Method	MSVOA_N HP Processing Method 82N112425W.M

STD. NAME	STD REF.#
Tune/Reschk Initial Calibration Stds	VP136451
CCC Internal Standard/PEM ICV/I.BLK Surrogate Standard MS/MSD Standard LCS Standard	VP136452,VP136453

Sr#	SampleID	ClientID	Data File Name	Date-Time	Comment	Operator	Status
1	BFB	BFB	VN088369.D	01 Dec 2025 08:49		JC\MD	Ok
2	VSTDCCC050	VSTDCCC050	VN088370.D	01 Dec 2025 09:58	pH#Lot#V12668	JC\MD	Ok,M
3	VN1201MBL01	VN1201MBL01	VN088371.D	01 Dec 2025 10:33		JC\MD	Ok
4	VN1201WBL01	VN1201WBL01	VN088372.D	01 Dec 2025 10:53		JC\MD	Ok
5	VN1201WBS01	VN1201WBS01	VN088373.D	01 Dec 2025 11:27		JC\MD	Ok,M
6	VN1201WBSD01	VN1201WBSD01	VN088374.D	01 Dec 2025 12:10		JC\MD	Ok,M
7	Q3715-02	MW6	VN088375.D	01 Dec 2025 12:30	vial A pH<2	JC\MD	Ok,M
8	Q3741-07	TB-1	VN088376.D	01 Dec 2025 12:51	vial A pH<2 TB	JC\MD	Ok
9	Q3716-01	MW5	VN088377.D	01 Dec 2025 13:11	vial A pH<2 Need 20X	JC\MD	Dilution
10	Q3716-03	MW4	VN088378.D	01 Dec 2025 13:32	Need Straight Run	JC\MD	Not Ok
11	Q3716-04	MW7	VN088379.D	01 Dec 2025 13:52	Need 200X	JC\MD	Dilution
12	Q3715-01	MW2	VN088380.D	01 Dec 2025 14:13	E flag of previous sample	JC\MD	ReRun
13	Q3716-02	MW3	VN088381.D	01 Dec 2025 14:33	vial A pH<2	JC\MD	Ok
14	Q3738-01	MW4	VN088382.D	01 Dec 2025 14:54	Run @ Lower Dilution	JC\MD	Not Ok
15	Q3738-03	Field	VN088383.D	01 Dec 2025 15:15	vial A pH<2 FB	JC\MD	Ok
16	Q3738-02	MW5	VN088384.D	01 Dec 2025 15:35	vial A pH<2	JC\MD	Ok,M
17	Q3745-01	STORAGE-BLANK-SO	VN088385.D	01 Dec 2025 15:56	vial A pH<2 Tics Contamination	JC\MD	ReRun
18	Q3745-02	STORAGE-BLANK-WA	VN088386.D	01 Dec 2025 16:16	vial A pH<2	JC\MD	Ok

Instrument ID: MSVOA_N

Daily Analysis Runlog For Sequence/QC Batch ID # VN120125

Review By	Semsettin Yesilyurt	Review On	12/4/2025 10:27:11 AM		
Supervise By	John Carlone	Supervise On	12/4/2025 10:36:07 AM		
SubDirectory	VN120125	HP Acquire Method	MSVOA_N	HP Processing Method	82N112425W.M
STD. NAME	STD REF.#				
Tune/Reschk Initial Calibration Stds	VP136451				
CCC Internal Standard/PEM ICV/I.BLK Surrogate Standard MS/MSD Standard LCS Standard	VP136452,VP136453				

19	Q3745-03	STORAGE-BLANK-WA	VN088387.D	01 Dec 2025 16:37	vial A pH<2	JC\MD	Ok
20	Q3745-04	STORAGE-BLANK-SAM	VN088388.D	01 Dec 2025 16:58	vial A pH<2	JC\MD	Ok
21	Q3743-03	TB	VN088389.D	01 Dec 2025 17:18	vial A pH<2 TB	JC\MD	Ok
22	Q3743-04	TB	VN088390.D	01 Dec 2025 17:39	vial A pH<2 TB	JC\MD	Ok
23	Q3743-05	TB	VN088391.D	01 Dec 2025 18:00	SIM Method on login	JC\MD	Not Ok
24	Q3743-01	TW-1125-1	VN088392.D	01 Dec 2025 18:20	vial A pH<2	JC\MD	Ok
25	Q3743-02	TW-1125-2	VN088393.D	01 Dec 2025 18:41	vial A pH<2	JC\MD	Ok
26	VSTDCCC050	VSTDCCC050EC	VN088394.D	01 Dec 2025 19:01		JC\MD	Ok,M

M : Manual Integration



SHIPPING DOCUMENTS

CLIENT INFORMATION

REPORT TO BE SENT TO:
COMPANY: Remington Vernick Engineers
ADDRESS: 2059 Springdale Road
CITY Cherry Hill STATE: NJ ZIP: 08003
ATTENTION: Kyle Carlson eRVE.com
PHONE: 609-682-3049 FAX:

CLIENT PROJECT INFORMATION

PROJECT NAME: Saddler Property
PROJECT NO.: 04001003 LOCATION: Camden, NJ
PROJECT MANAGER: Kyle Carlson
e-mail: Kyle.Carlson@RVE.com
PHONE: 609-682-3049 FAX:

CLIENT BILLING INFORMATION

BILL TO: AP Invoices eRVE.com PO#: 04001003
ADDRESS: 2059 Springdale Road
CITY Cherry Hill STATE: NJ ZIP: 08003
ATTENTION: Kyle Carlson PHONE: 609-682-3049

ANALYSIS

DATA TURNAROUND INFORMATION

FAX (RUSH) _____ DAYS*
HARDCOPY (DATA PACKAGE): Standard 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 NJ HazSite

1: PCB
2: Pesticide TCL
3: SVOC-TCL-BNA20
4: VOC-TCL-VOA+ID
5: VOC SIM
6: Heavy Metals TCLICP
7: Cyanide
8:
9:

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		E	E	E	A	A	B	D		
1.	TW - 1125-1	GW		X	11/26/25	915	9	1	1	1	2	2	1	1		
2.	TW - 1125-2	GW		X	11/26/25	1000	9	1	1	1	2	2	1	1		
3.	TB				11/26/25	1000	2				2					
4.	TB				11/26/25	1000	2				2					
5.	TB				11/26/25	1000	2					2				
6.																
7.																
8.																
9.																
10.																

SAMPLE CUSTODY MUST BE DOCUMENTED BELOW EACH TIME SAMPLES CHANGE POSSESSION INCLUDING COURIER DELIVERY

RELINQUISHED BY SAMPLER: 1. <u>Sutton Zook eRVE.com</u>	DATE/TIME: <u>11/26/25</u>	RECEIVED BY: <u>[Signature]</u> <u>11-26-25</u>	Conditions of bottles or coolers at receipt: ☐ COMPLIANT ☐ NON COMPLIANT ☐ COOLER TEMP <u>2.8</u> °C
RELINQUISHED BY SAMPLER: 2.	DATE/TIME:	RECEIVED BY: 2.	Comments:
RELINQUISHED BY SAMPLER: 3.	DATE/TIME: <u>11-26-25</u>	RECEIVED BY: 3.	Page <u>1</u> of <u>1</u> CLIENT: ☐ Hand Delivered ☐ Other

Shipment Complete
☐ YES ☐ NO

From: Kyle Carlson <Kyle.Carlson@rve.com>
Sent: Monday, December 01, 2025 1:56 PM
Subject: Re: Sadler Project - Low Volume Sample #1

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Secured by Check Point

Hi Deepak - confirmed to proceed with the SVOC full scan without the SIM due to the low volume provided. Thank you.

Regards,

Kyle Carlson, LSRP
Environmental Project Manager
Remington & Vernick Engineers
2059 Springdale Road
Cherry Hill, NJ 08003
Cell: (609) 682-3049
Kyle.Carlson@rve.com



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From: Deepak Parmar <Deepak.Parmar@alliancetg.com>
Sent: Monday, December 1, 2025 1:50 PM
To: Kyle Carlson <Kyle.Carlson@rve.com>
Cc: Justin Zarzecki <Justin.Zarzecki@rve.com>
Subject: Re: Sadler Project - Low Volume Sample #1

Hi Kyle,

We haven't received extra volume to perform SVOC-SIM analysis in order to meet the NJ GWQS limits. Therefore, we would like to confirm that lab will proceed with only SVOC full scan analysis with the received sample volume.

Thanks & Regards,





Deepak Parmar
Sr. Project Manager
An Alliance Technical Group Company
Main: 908-789-8900
Direct: 908-728-3154
Address: 284 Sheffield St, Ste 1, Mountainside, NJ 07092
www.alliancetg.com



From: Kyle Carlson <Kyle.Carlson@rve.com>
Sent: Monday, December 1, 2025 10:28 AM
To: Deepak Parmar <deepak.parmar@alliancetg.com>
Cc: Justin Zarzecki <Justin.Zarzecki@rve.com>
Subject: Sadler Project - Low Volume Sample #1

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Secured by Check Point

Hi Deepak - confirming run of SVOCs on sample #1 (TW-1125-1) without the PCBs and Pesticides analyses due to the low-volume provided. Please let me know if I can provide any additional info - thank you.

Regards,

Kyle Carlson, LSRP
Environmental Project Manager
Remington & Vernick Engineers
2059 Springdale Road
Cherry Hill, NJ 08003
Cell: (609) 682-3049
Kyle.Carlson@rve.com



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From: Kyle Carlson <Kyle.Carlson@rve.com>
Sent: Monday, December 01, 2025 10:28 AM
Subject: Sadler Project - Low Volume Sample #1

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Secured by Check Point

Hi Deepak - confirming run of SVOCs on sample #1 (TW-1125-1) without the PCBs and Pesticides analyses due to the low-volume provided. Please let me know if I can provide any additional info - thank you.

Regards,

Kyle Carlson, LSRP
Environmental Project Manager
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2059 Springdale Road
Cherry Hill, NJ 08003
Cell: (609) 682-3049
Kyle.Carlson@rve.com



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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



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

LOGIN REPORT/SAMPLE TRANSFER

Order ID : Q3743 REMI01
Client Name : Remington & Vernick
Client Contact : Justin Zarzecki
Invoice Name : Remington & Vernick
Invoice Contact : Justin Zarzecki

Order Date : 12/1/2025 9:45:00 AM
Project Name : *Saddler* Sadler Property
Receive DateTime : 11/26/2025 6:20:00 PM
Purchase Order :

Project Mgr : Deepak
Report Type : Level 1 *Regular*
EDD Type : ~~EXCEL NUCLEANUP~~
Hard Copy Date : *HT Reducel*
Date Signoff : 12/1/2025 1:00:29 PM

LAB ID	CLIENT ID	MATRIX	SAMPLE DATE	SAMPLE TIME	TEST	TEST GROUP	METHOD	FAX DATE	DUE DATES
Q3743-01	TW-1125-1	Water	11/26/2025	09:15	VOC-SIM		SFAM_VOCSIM	10 Bus. Days	
					VOC-TCLVOA-10	TCL+30/TAL	8260-Low	10 Bus. Days	
Q3743-02	TW-1125-2	Water	11/26/2025	10:00	VOC-SIM		SFAM_VOCSIM	10 Bus. Days	
					VOC-TCLVOA-10	TCL+30/TAL	8260-Low	10 Bus. Days	
Q3743-03	TB	Water	11/26/2025	10:00	VOC-TCLVOA-10	TCL+30/TAL	8260-Low	10 Bus. Days	
Q3743-04	TB	Water	11/26/2025	10:00	VOC-TCLVOA-10	TCL+30/TAL	8260-Low	10 Bus. Days	
Q3743-05	TB	Water	11/26/2025	10:00	VOC-SIM		SFAM_VOCSIM	10 Bus. Days	



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

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Client Contact : Justin Zarzecki
Invoice Name : Remington & Vernick
Invoice Contact : Justin Zarzecki

Order Date : 12/1/2025 9:45:00 AM
Project Name : *Sadler* Sadler Property
Receive DateTime : 11/26/2025 6:20:00 PM
Purchase Order :

Project Mgr : Deepak
Report Type : Level 1 *Regular*
EDD Type : EXCEL NUCLEONUP
Hard Copy Date : *MS Redwood*
Date Signoff : 12/1/2025 1:00:29 PM

LAB ID	CLIENT ID	MATRIX	SAMPLE DATE	SAMPLE TIME	TEST	TEST GROUP	METHOD	FAX DATE	DUE DATES
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*Spilled in ref
left 04*

Relinquished By : *[Signature]*

Date / Time : *12/1/25 14:21*

Received By : *[Signature]*

Date / Time : *12/1/25 14:30*

Storage Area : VOA Refrigerator Room