

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN042924\
Data File : VN081908.D
Acq On : 29 Apr 2024 12:44
Operator : JC\MD
Sample : P2292-01
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
240418050-02

Integration Parameters: RTEINT3.P

Integrator: RTE

Smoothing : OFF

Sampling : 1

Start Thrs: 0.2

Stop Thrs : 0

Filtering: 5

Min Area: 0 % 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\624N042624W.M

Title : METHOD 624 VOLATILE ORGANIC ANALYSIS

Signal : TIC: VN081908.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.283	47	56	61	rBV	3819806	6905688	100.00%	65.394%
2	4.424	408	420	431	rBV	59137	173339	2.51%	1.641%
3	7.818	985	997	1009	rBV2	127190	342792	4.96%	3.246%
4	8.577	1118	1126	1138	rBV	136848	310002	4.49%	2.936%
5	9.106	1206	1216	1227	rBV	297253	624098	9.04%	5.910%
6	10.571	1456	1465	1473	rBV	466483	875053	12.67%	8.286%
7	11.865	1677	1685	1699	rVB	398726	712513	10.32%	6.747%
8	12.853	1844	1853	1861	rBV	288249	504314	7.30%	4.776%
9	14.247	2083	2090	2095	rBV	70020	112338	1.63%	1.064%

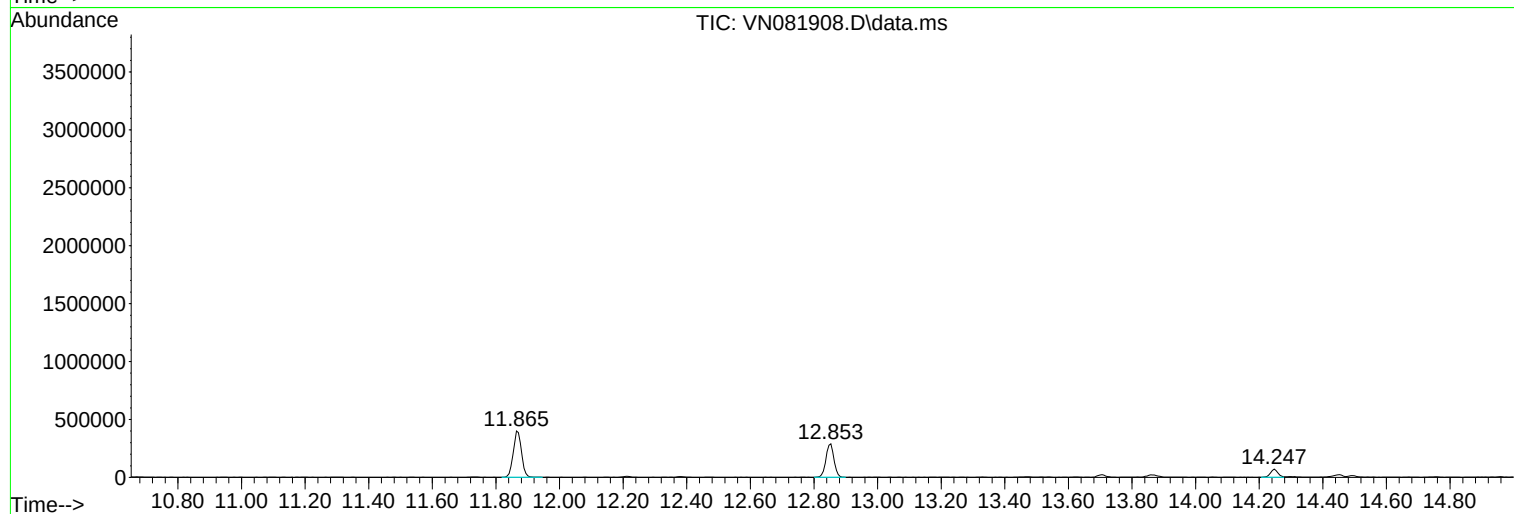
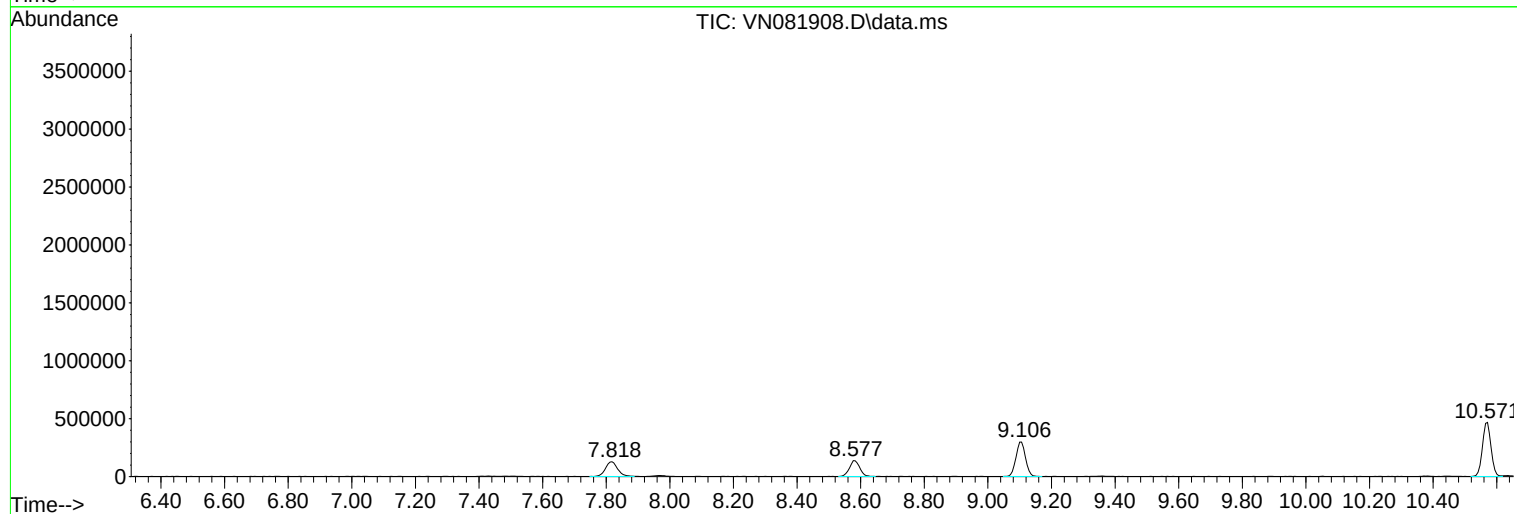
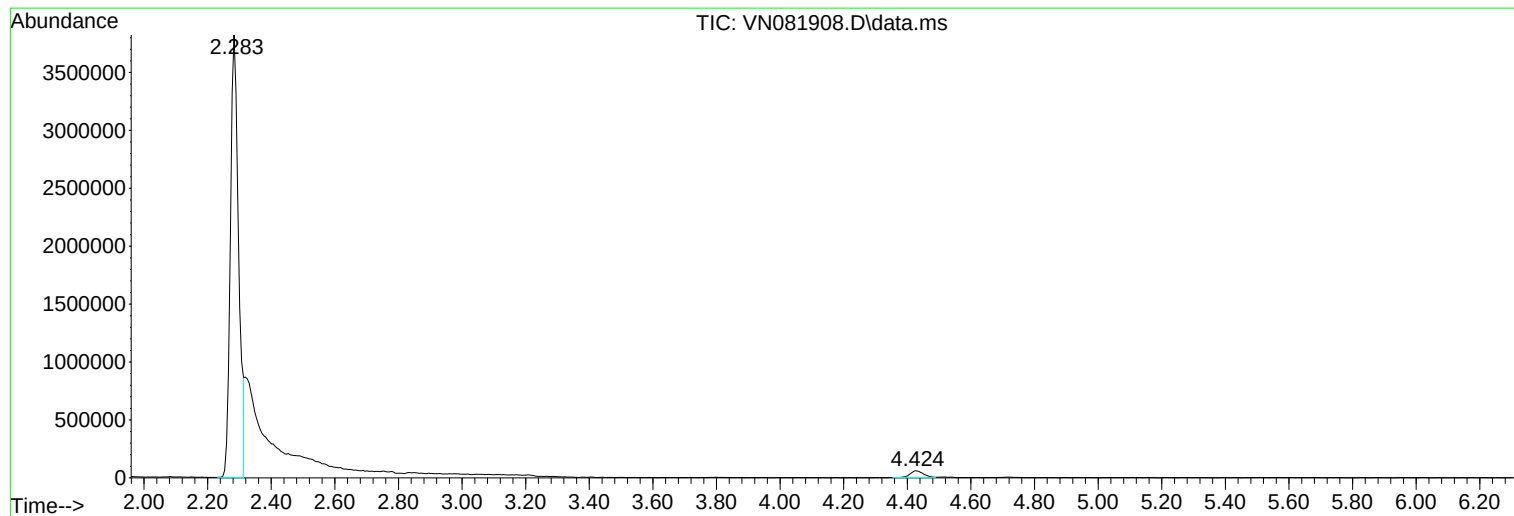
Sum of corrected areas: 10560137

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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\624N042624W.M
Quant Title : METHOD 624 VOLATILE ORGANIC ANALYSIS

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



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ClientSampleId :
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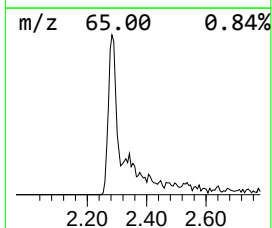
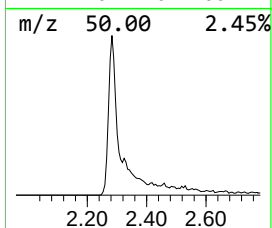
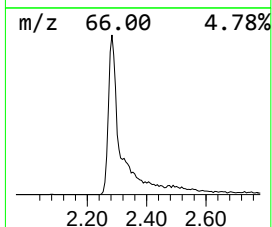
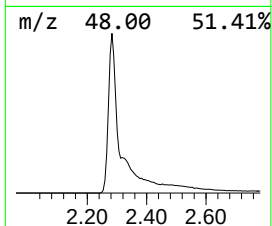
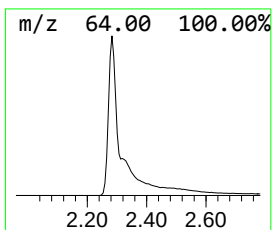
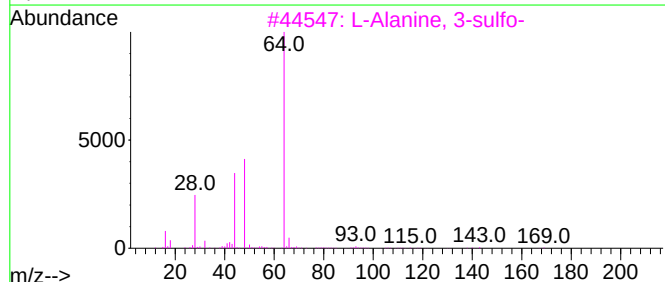
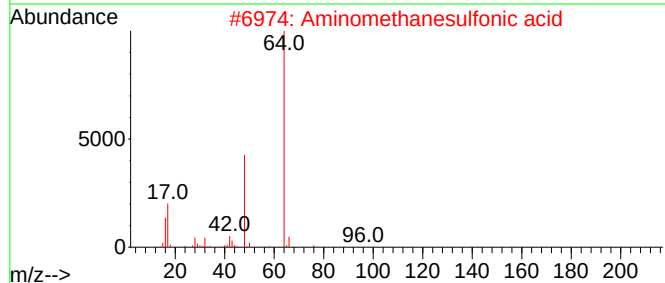
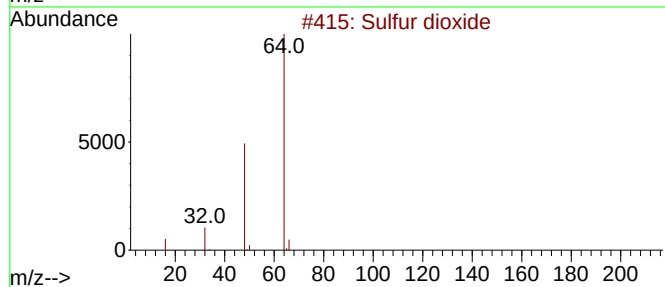
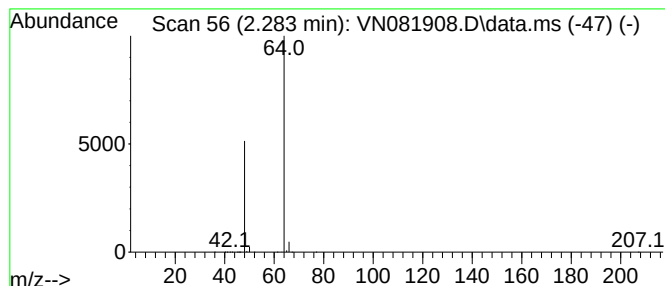
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\624N042624W.M
Quant Title : METHOD 624 VOLATILE ORGANIC ANALYSIS

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

Peak Number 1 Sulfur dioxide Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.283	604.36 ug/l	6905690	Bromochloromethane	7.818

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Sulfur dioxide	64	O2S	007446-09-5	83
2			Aminomethanesulfonic acid	111	CH5NO3S	013881-91-9	74
3			L-Alanine, 3-sulfo-	169	C3H7NO5S	000498-40-8	9
4			Aminomethanesulfonic acid	111	CH5NO3S	013881-91-9	9
5			Ethene, 1,1-difluoro-	64	C2H2F2	000075-38-7	4



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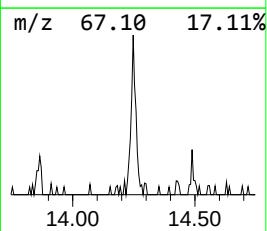
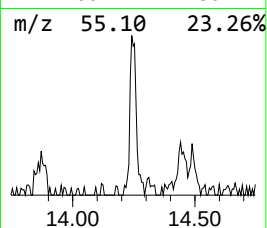
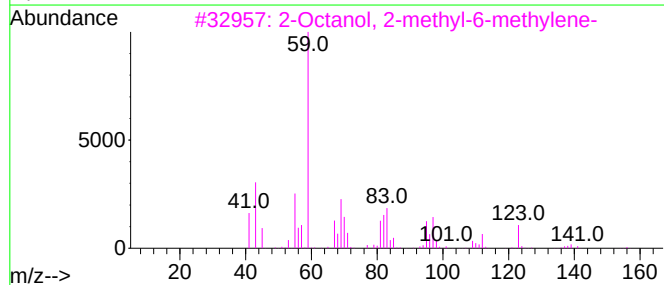
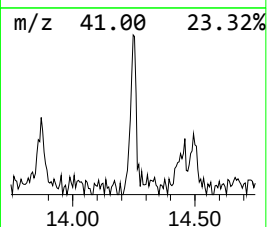
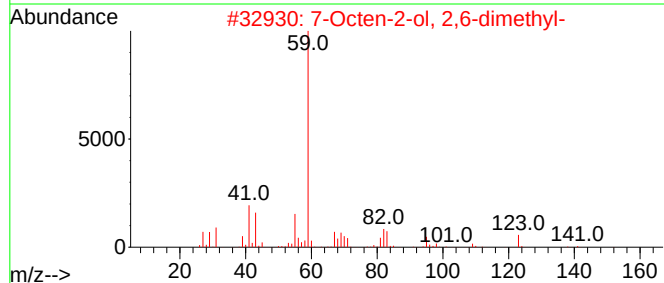
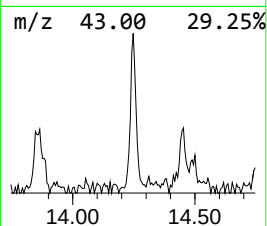
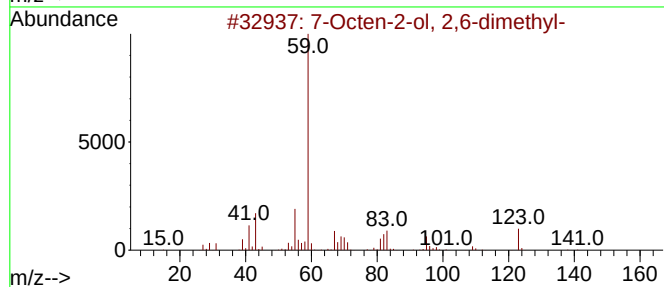
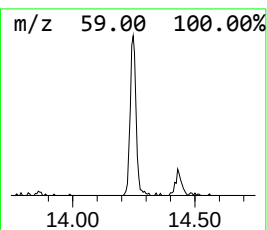
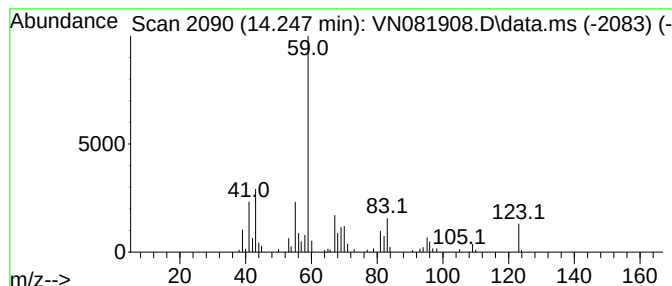
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TIC Integration Parameters: LSCINT.P

Peak Number 2 7-Octen-2-ol, 2,6-dimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.247	4.73 ug/l	112338	Chlorobenzene-d5	11.865

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			7-Octen-2-ol, 2,6-dimethyl-	156	C10H20O	018479-58-8	83
2			7-Octen-2-ol, 2,6-dimethyl-	156	C10H20O	018479-58-8	64
3			2-Octanol, 2-methyl-6-methylene-	156	C10H20O	018479-59-9	64
4			1,7-Octanediol, 3,7-dimethyl-	174	C10H22O2	000107-74-4	56
5			1,7-Octanediol, 3,7-dimethyl-	174	C10H22O2	000107-74-4	50



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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Sulfur dioxide	2.283	604.4	ug/l	6905690	1	7.818	342792	30.0
7-Octen-2-ol, 2...	14.247	4.7	ug/l	112338	3	11.865	712513	30.0