

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN022422\
Data File : VN071065.D
Acq On : 24 Feb 2022 19:08
Operator : JC\MD
Sample : N1638-02
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
40473

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\82N021522W.M

Title : SW846 8260

Signal : TIC: VN071065.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.540	91	94	125	rBV2	47322121	193124317	15.89%	12.236%
2	4.157	275	369	374	rBV4	46600618	1215202895	100.00%	76.995%
3	4.340	389	400	426	rVB	3580917	13503588	1.11%	0.856%
4	5.634	604	620	640	rBV	4815969	16757443	1.38%	1.062%
5	8.257	1054	1066	1085	rBV	12114314	28794652	2.37%	1.824%
6	8.451	1085	1099	1111	rBV	19006311	44696289	3.68%	2.832%
7	10.498	1441	1447	1463	rVB	24228202	44583865	3.67%	2.825%
8	13.763	1996	2002	2010	rBV	13938697	21629445	1.78%	1.370%

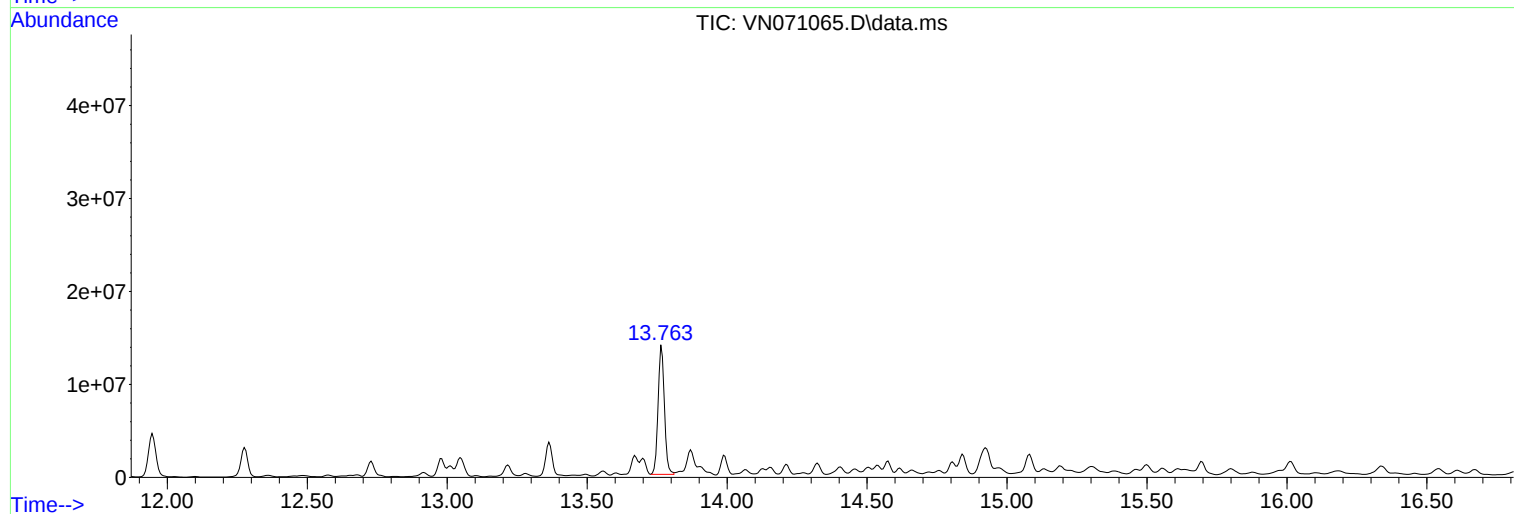
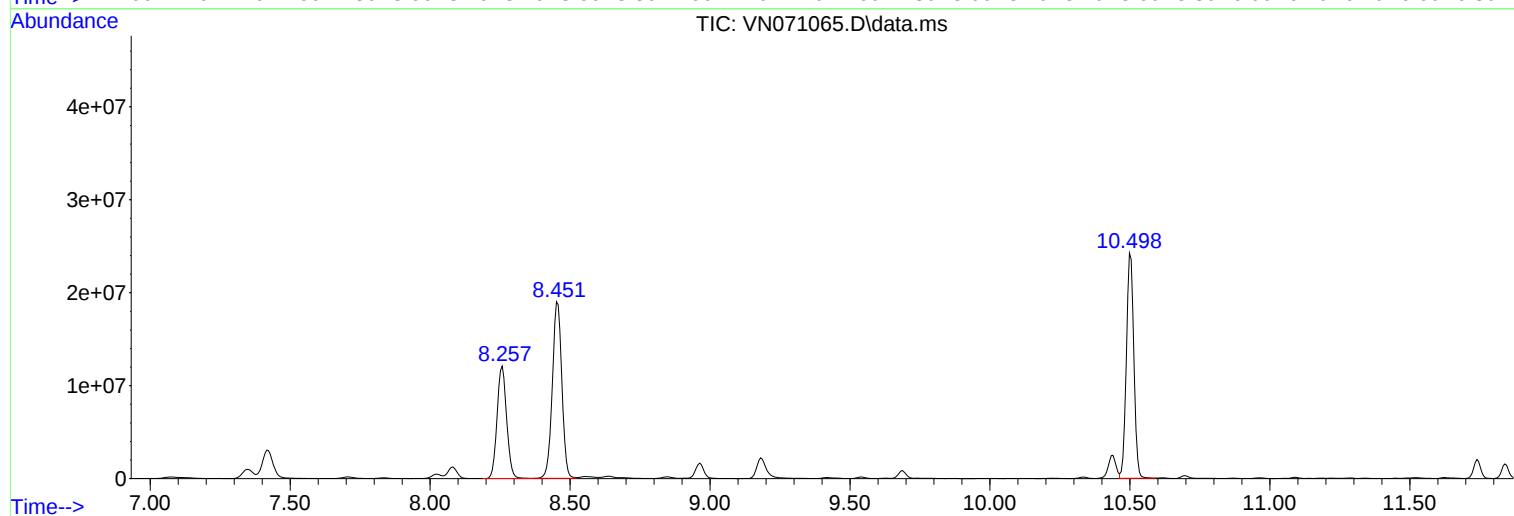
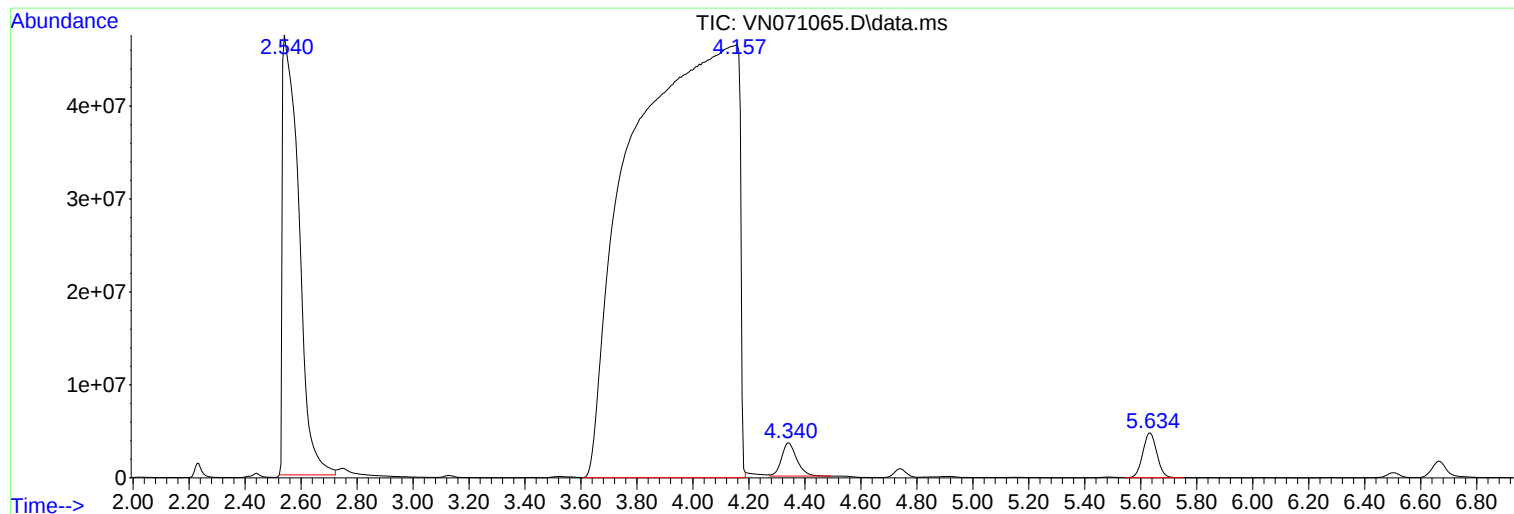
Sum of corrected areas: 1578292494

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Quant Title : SW846 8260

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



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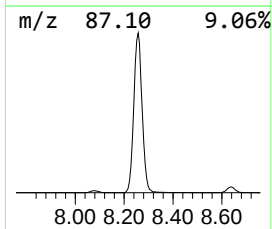
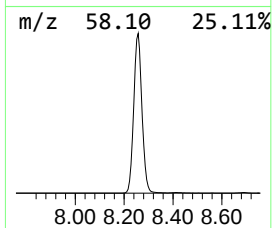
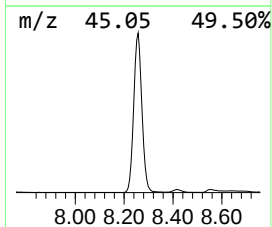
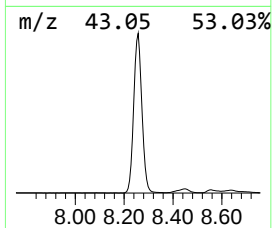
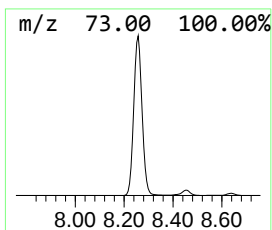
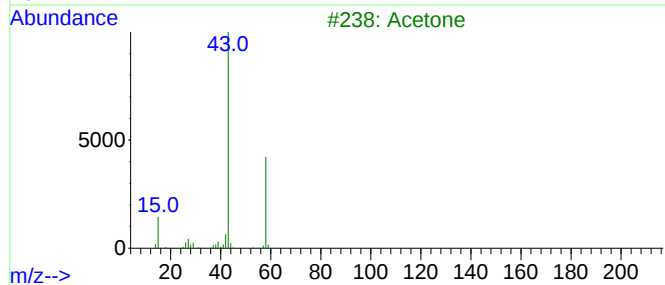
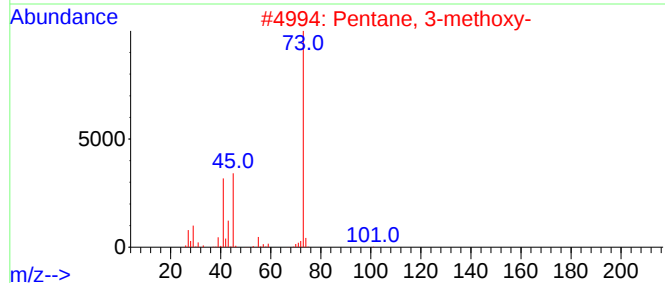
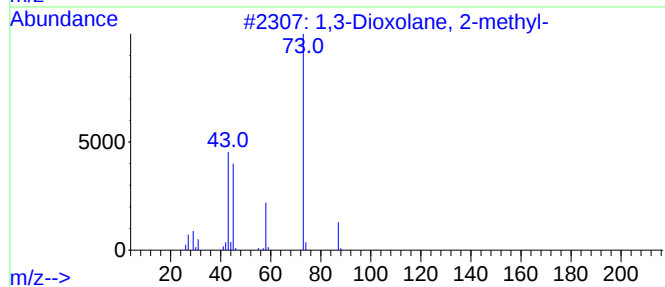
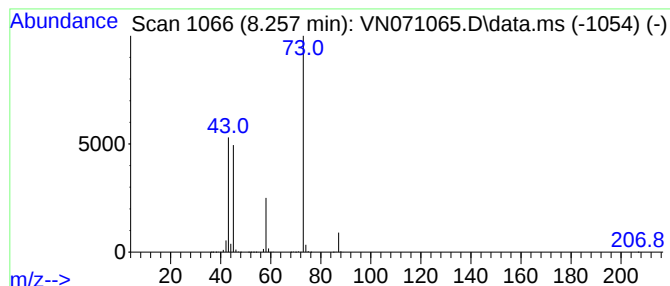
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N021522W.M
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TIC Library : C:\Database\NIST0.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 1,3-Dioxolane, 2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.257	50.00 ug/l	28794700	Pentafluorobenzene	8.081

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	91
2			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	28
3			Acetone	58	C3H6O	000067-64-1	9
4			Silane, tetramethyl-	88	C4H12Si	000075-76-3	9
5			2-Acetamido-N-methylacetamide	130	C5H10N2O2	007606-79-3	9



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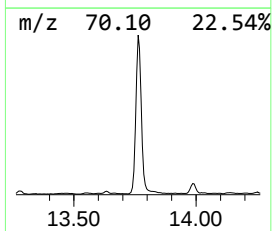
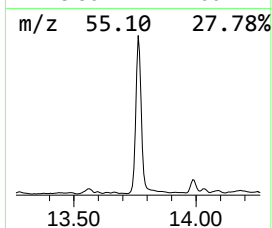
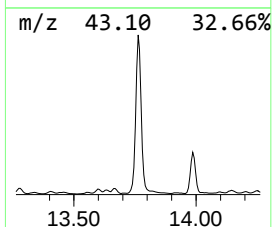
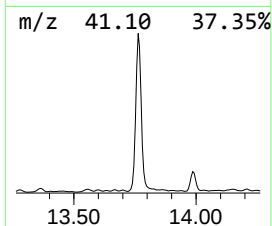
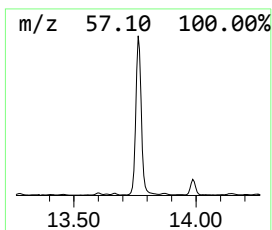
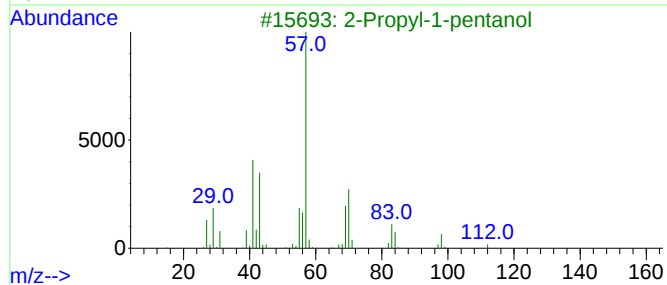
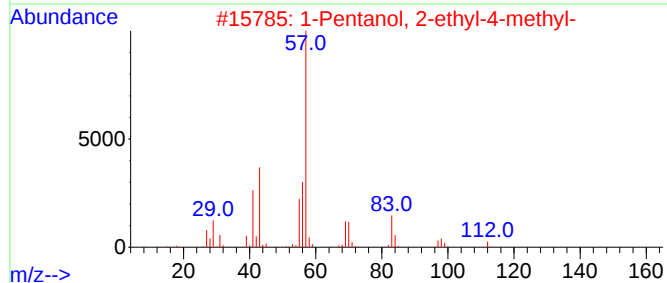
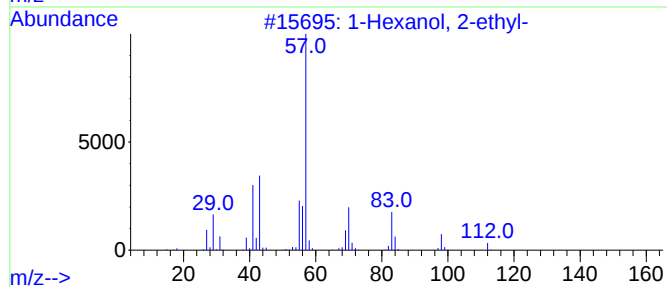
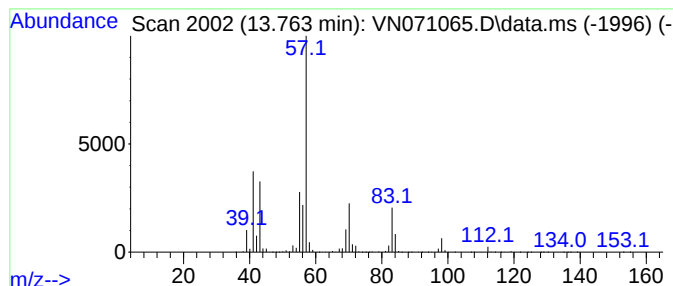
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TIC Integration Parameters: LSCINT.P

Peak Number 4 1-Hexanol, 2-ethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.763	50.00 ug/l	21629400	1,4-Dichlorobenzene-d4	13.669

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	83
2			1-Pentanol, 2-ethyl-4-methyl-	130	C8H18O	000106-67-2	56
3			2-Propyl-1-pentanol	130	C8H18O	058175-57-8	53
4			Pentadecafluorooctanoic acid, 2-...	526	C16H17F15O2	1000406-80-9	47
5			1-Decene, 2,4-dimethyl-	168	C12H24	055170-80-4	47



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
1,3-Dioxolane, ...	8.257	50.0	ug/l	28794700	1	8.081	28794700	50.0
1-Hexanol, 2-et...	13.763	50.0	ug/l	21629400	4	13.669	21629400	50.0