

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN091521\
 Data File : VN068491.D
 Acq On : 15 Sep 2021 18:15
 Operator : JC/MD
 Sample : M3759-09
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 31206

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\82N090721W.M

Title : SW846 8260

Signal : TIC: VN068491.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.140	753	802	846	rBV2	27374421	68061600	100.00%	81.185%
2	6.581	1683	1712	1765	rBV2	893159	2901406	4.26%	3.461%
3	8.024	2228	2250	2261	rBV	369774	866405	1.27%	1.033%
4	8.088	2262	2274	2309	rVB	720264	1629638	2.39%	1.944%
5	8.440	2384	2405	2430	rBV	335625	770314	1.13%	0.919%
6	8.971	2585	2603	2626	rBV	870685	1803607	2.65%	2.151%
7	10.446	3131	3153	3189	rBV	1305908	2456925	3.61%	2.931%
8	11.749	3614	3639	3667	rBV	1136653	2117236	3.11%	2.525%
9	12.736	3990	4007	4026	rBV	863556	1504847	2.21%	1.795%
10	13.677	4341	4358	4378	rBV	1000639	1723481	2.53%	2.056%

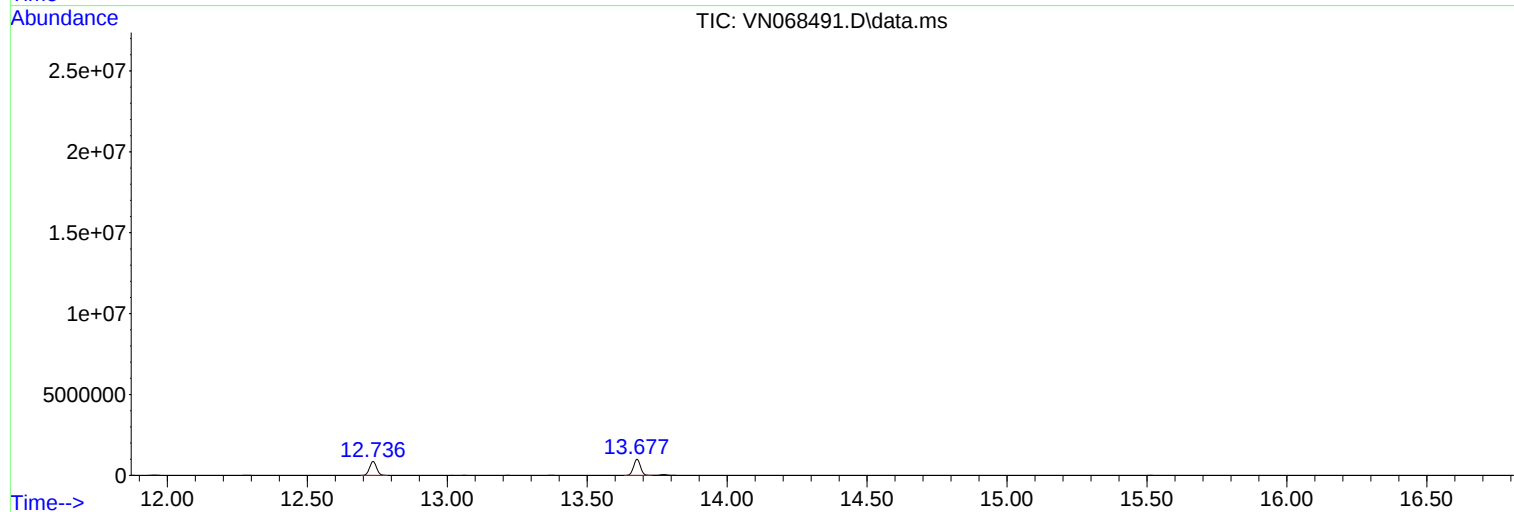
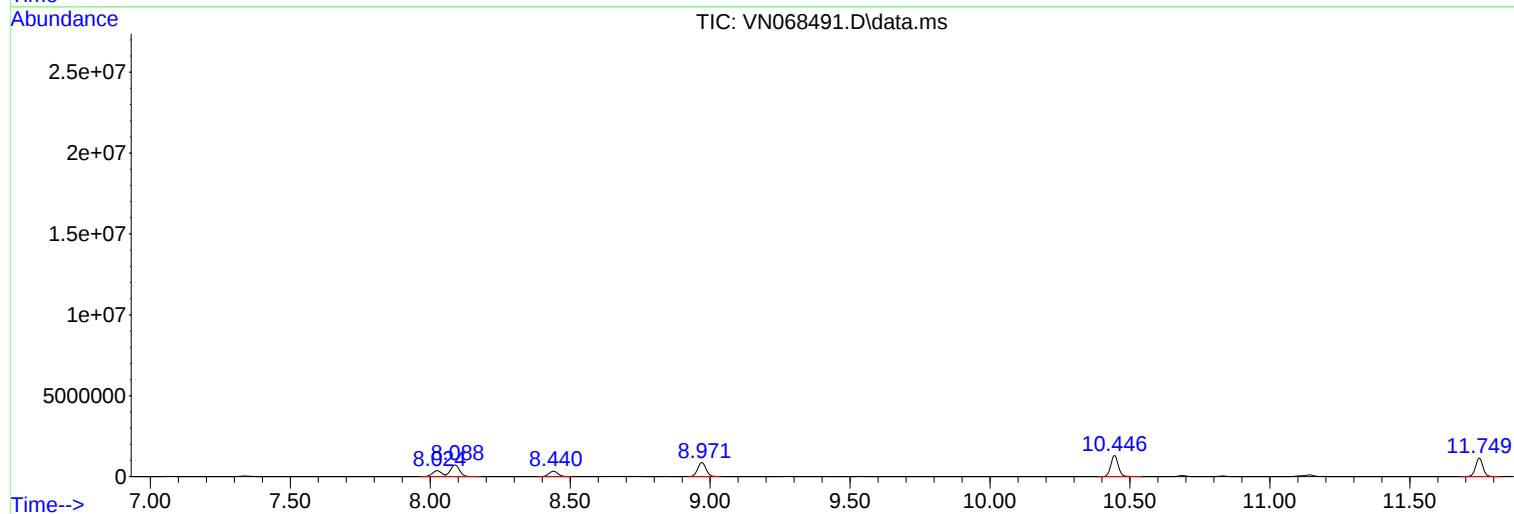
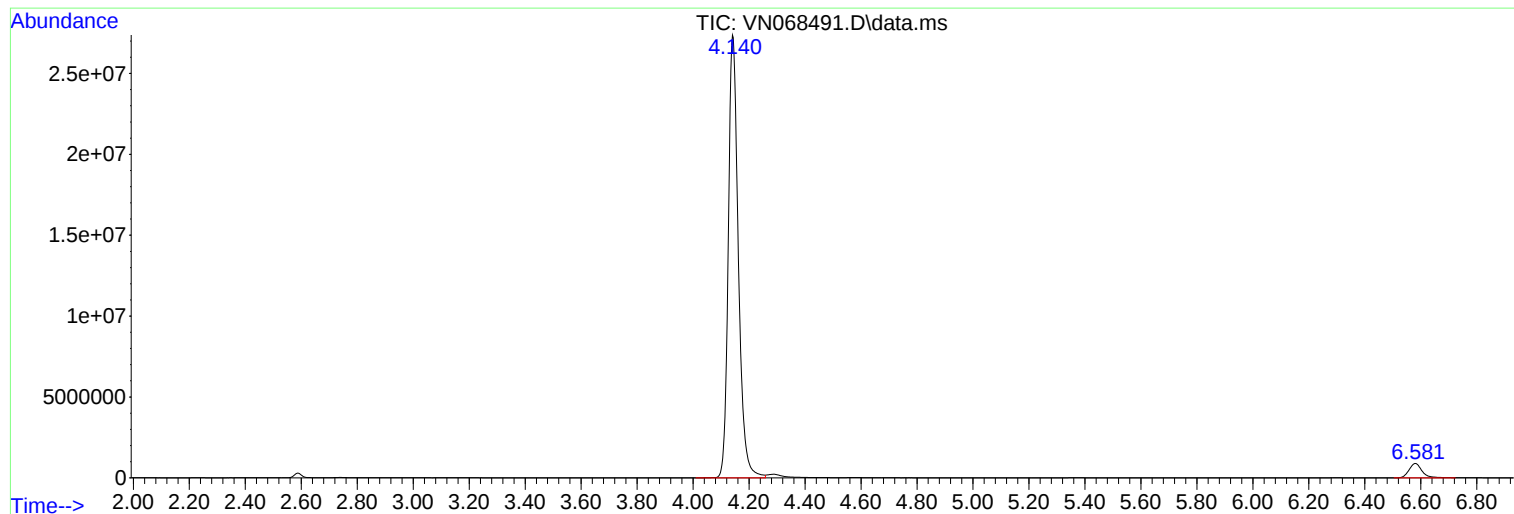
Sum of corrected areas: 83835459

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Quant Title : SW846 8260

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



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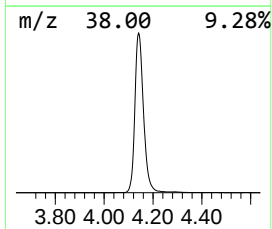
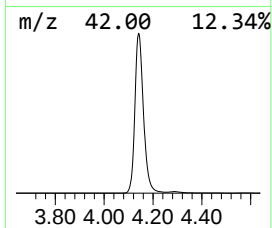
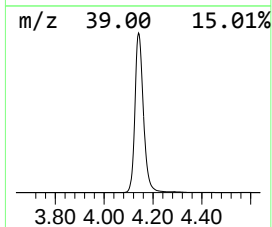
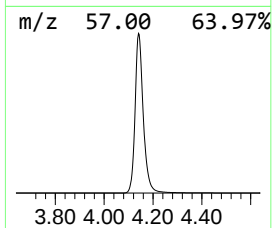
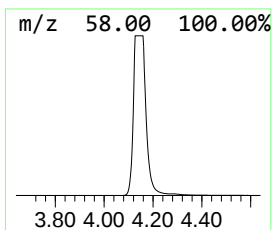
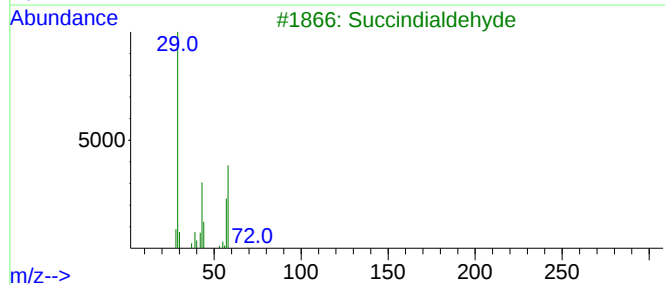
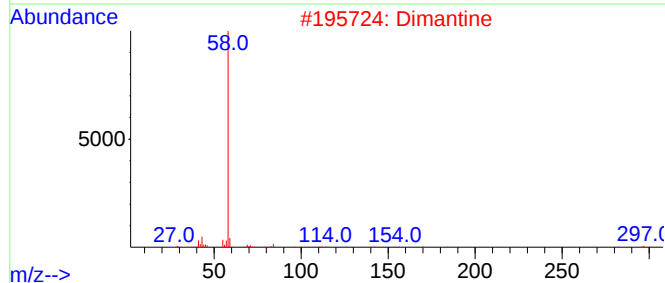
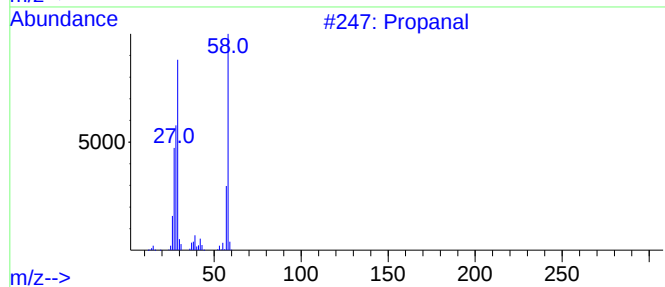
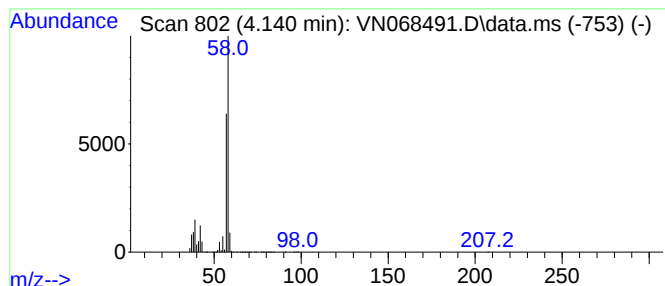
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Peak Number 1 unknown4.140 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.140	2088.24 ug/l	68061600	Pentafluorobenzene	8.088

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propanal	58	C3H6O	000123-38-6	9
2			Dimantine	297	C20H43N	000124-28-7	9
3			Succindialdehyde	86	C4H6O2	000638-37-9	9
4			Oxirane, methyl-, (S)-	58	C3H6O	016088-62-3	7
5			Methylamine, N,N-dimethyl-	59	C3H9N	000075-50-3	4



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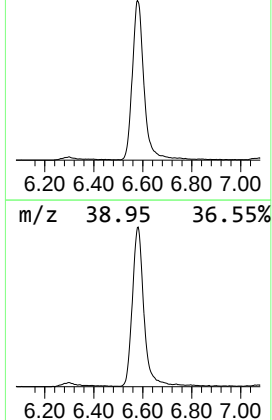
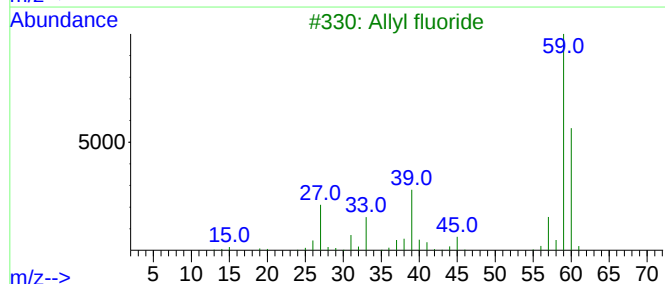
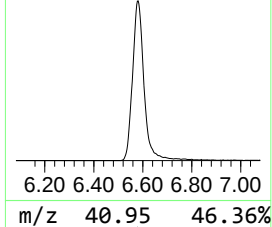
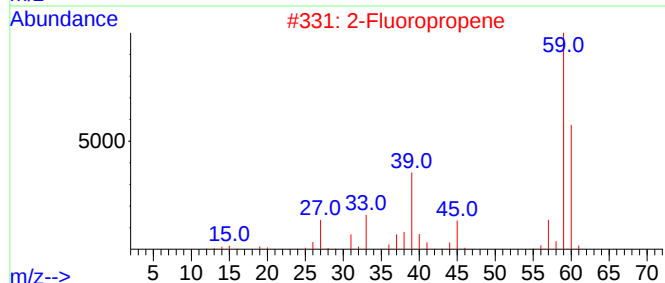
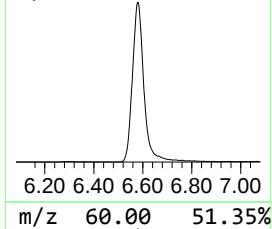
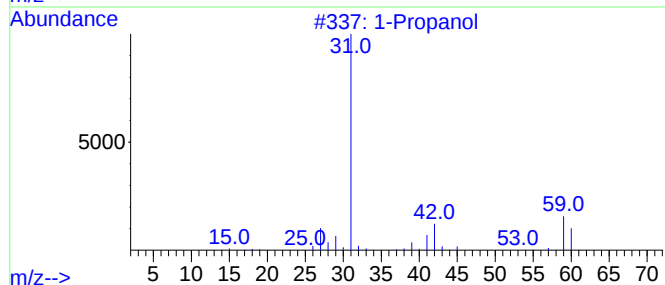
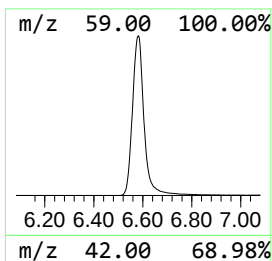
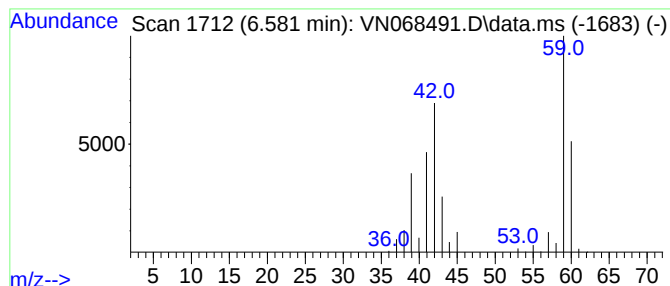
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Peak Number 2 1-Propanol Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.581	89.02 ug/l	2901410	Pentafluorobenzene	8.088

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Propanol	60	C3H8O	000071-23-8	72
2			2-Fluoropropene	60	C3H5F	001184-60-7	64
3			Allyl fluoride	60	C3H5F	000818-92-8	53
4			Cyclopropane	42	C3H6	000075-19-4	9
5			Acetaldoxime	59	C2H5NO	000107-29-9	7



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown4.140	4.140	2088.2	ug/l	68061600	1	8.088	1629640	50.0
1-Propanol	6.581	89.0	ug/l	2901410	1	8.088	1629640	50.0