

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN031221\
 Data File : VN066192.D
 Acq On : 12 Mar 2021 19:21
 Operator : JC/MD
 Sample : M1647-20
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 COMP4-41735-30780

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\82N031221W.M
 Title : SW846 8260

Signal : TIC: VN066192.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.621	698	720	748	rBV	14571155	42299375	100.00%	65.600%
2	3.739	750	764	803	rVB2	1769252	4676505	11.06%	7.253%
3	5.957	1565	1591	1647	rBV	1478952	4721874	11.16%	7.323%
4	7.512	2147	2171	2184	rBV3	314575	781756	1.85%	1.212%
5	7.587	2184	2199	2223	rVB2	593144	1395054	3.30%	2.164%
6	7.952	2310	2335	2359	rBV3	487376	1226673	2.90%	1.902%
7	8.515	2528	2545	2573	rVB2	877844	1836481	4.34%	2.848%
8	10.028	3091	3109	3125	rBV	1394468	2515127	5.95%	3.901%
9	11.350	3582	3602	3627	rBV	1145436	2061745	4.87%	3.197%
10	12.345	3948	3973	4002	rBV2	788661	1548425	3.66%	2.401%
11	13.286	4310	4324	4354	rBV	767499	1417258	3.35%	2.198%

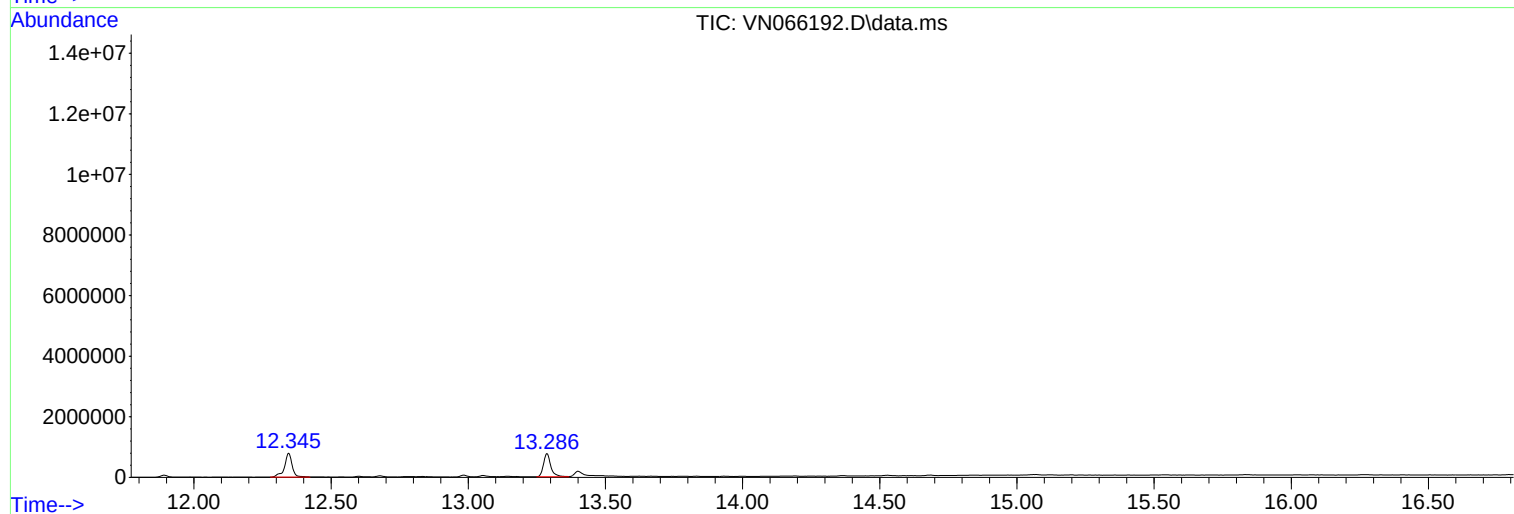
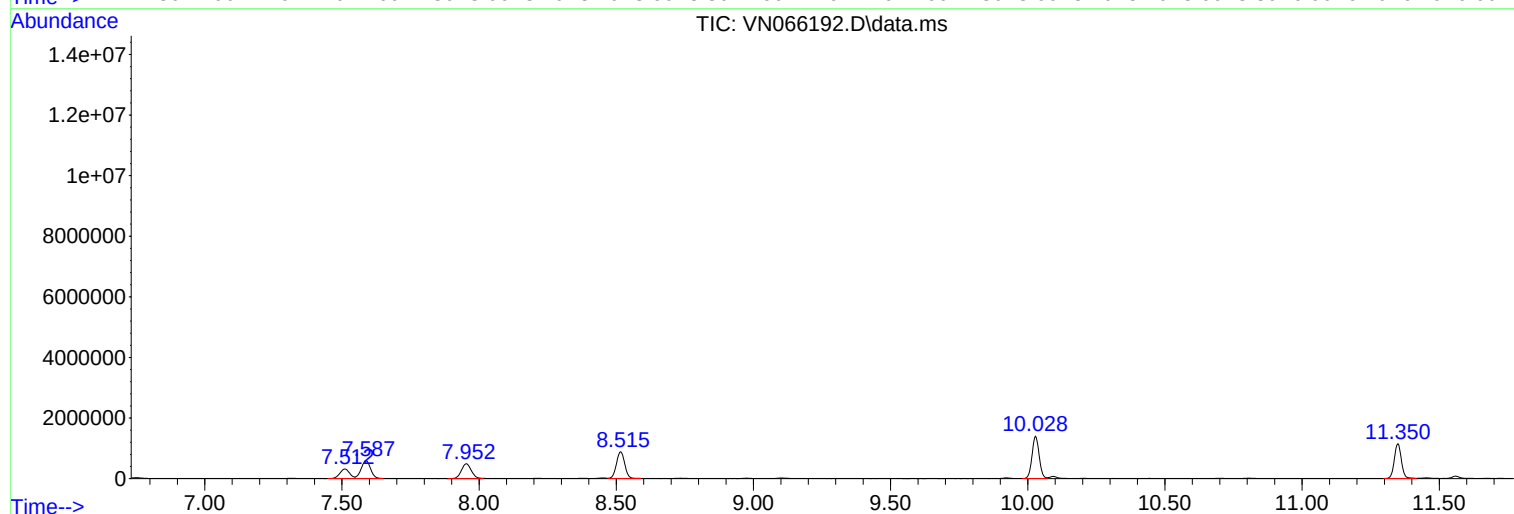
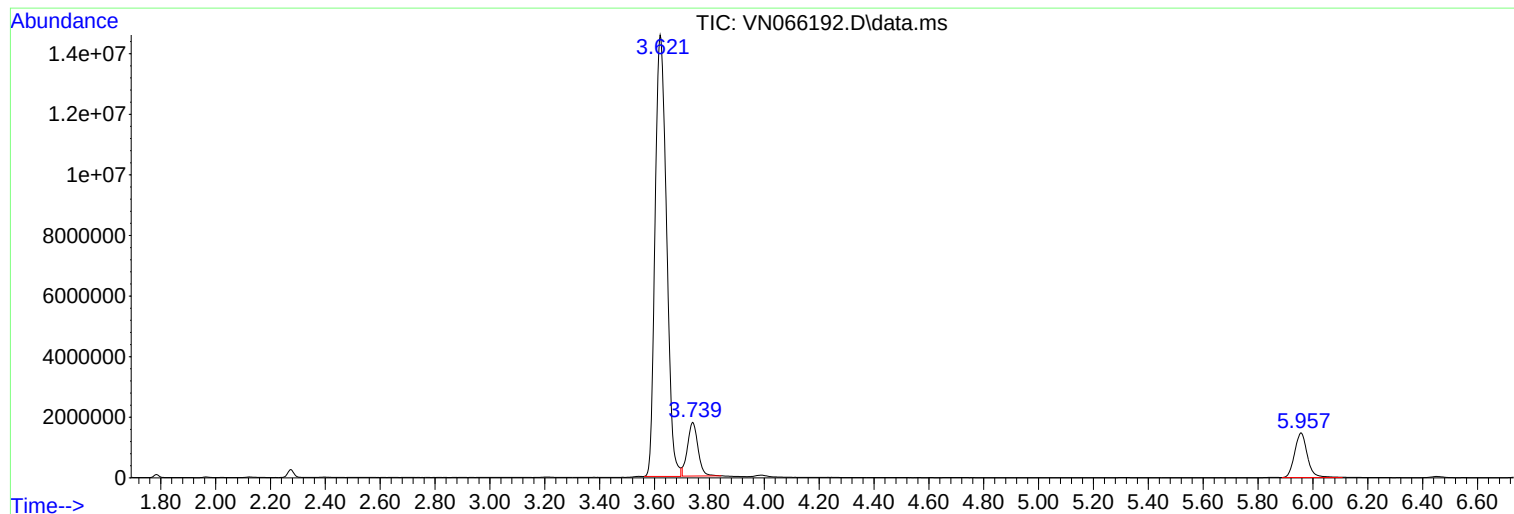
Sum of corrected areas: 64480273

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Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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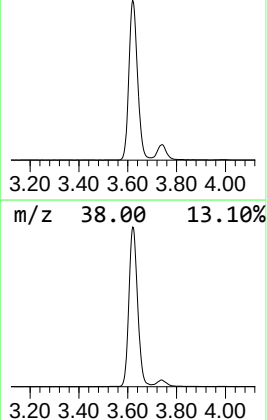
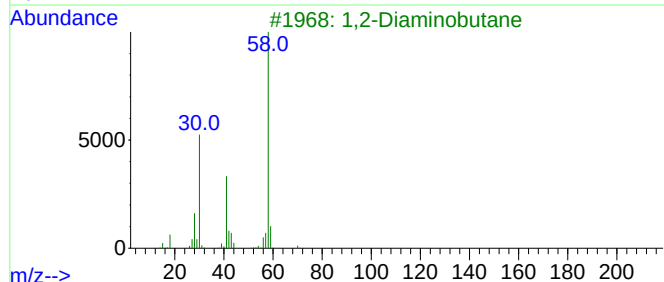
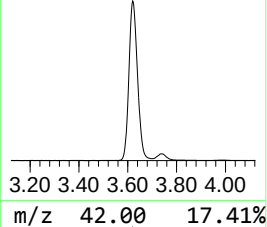
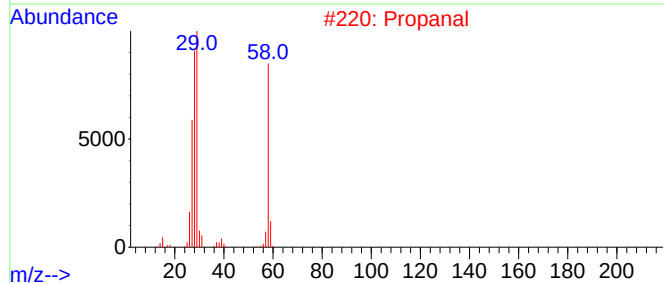
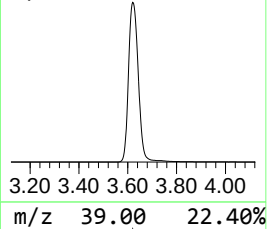
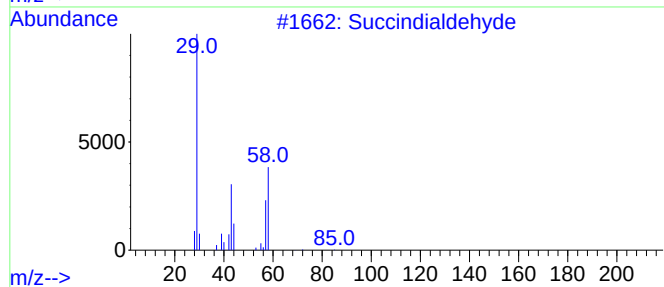
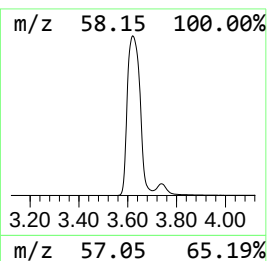
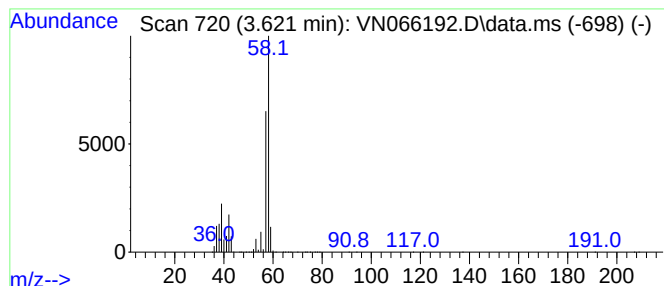
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Peak Number 1 unknown3.621 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.621	1516.05 ug/l	42299400	Pentafluorobenzene	7.587

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Succindialdehyde	86	C4H6O2	000638-37-9	40
2			Propanal	58	C3H6O	000123-38-6	9
3			1,2-Diaminobutane	88	C4H12N2	1000343-51-4	9
4			Methylamine, N,N-dimethyl-	59	C3H9N	000075-50-3	7
5			Oxirane, methyl-, (S)-	58	C3H6O	016088-62-3	5



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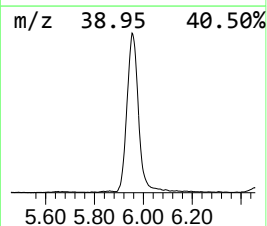
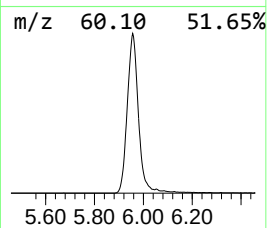
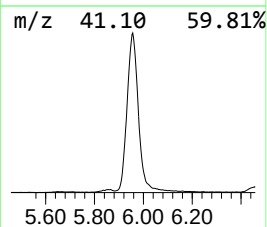
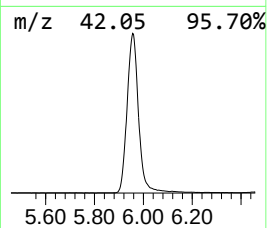
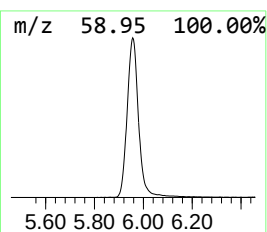
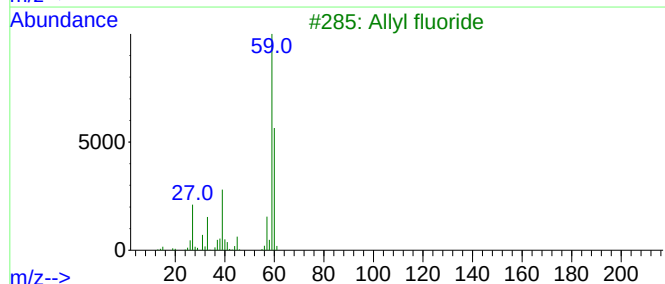
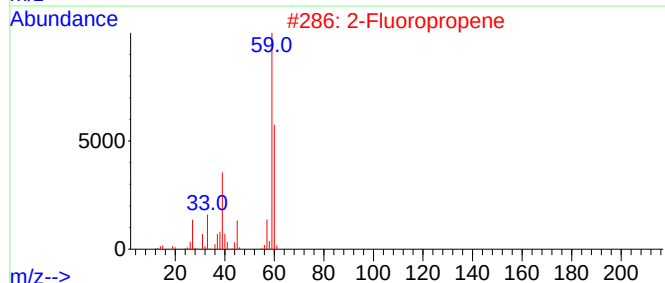
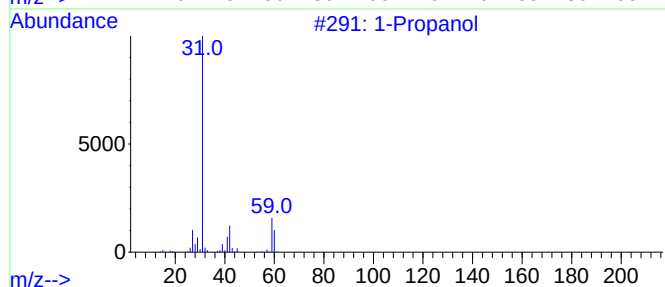
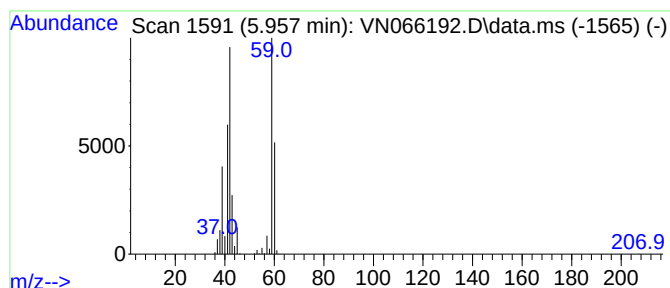
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Peak Number 2 1-Propanol Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.957	169.24 ug/l	4721870	Pentafluorobenzene	7.587

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Propanol	60	C3H8O	000071-23-8	64
2			2-Fluoropropene	60	C3H5F	001184-60-7	50
3			Allyl fluoride	60	C3H5F	000818-92-8	32
4			Propene	42	C3H6	000115-07-1	9
5			Cyclopropane	42	C3H6	000075-19-4	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown3.621	3.621	1516.1	ug/l	42299400	1	7.587	1395050	50.0
1-Propanol	5.957	169.2	ug/l	4721870	1	7.587	1395050	50.0