

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX031822\
 Data File : VX027533.D
 Acq On : 18 Mar 2022 17:17
 Operator : JC/MD
 Sample : N1998-04
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-130

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_X\Method\82X031722W.M

Title : SW846 8260

Signal : TIC: VX027533.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.965	297	308	323	rBV	2610335	5692544	25.37%	14.533%
2	3.117	323	333	366	rVB	9512105	22437233	100.00%	57.283%
3	4.318	519	530	545	rBV	91907	275445	1.23%	0.703%
4	5.391	694	706	722	rBV2	213473	622686	2.78%	1.590%
5	5.556	722	733	750	rVB	347142	988185	4.40%	2.523%
6	5.964	789	800	815	rBV	208819	564018	2.51%	1.440%
7	6.330	851	860	875	rVB	407197	1131091	5.04%	2.888%
8	6.763	920	931	950	rBV	536401	1289807	5.75%	3.293%
9	8.653	1234	1241	1257	rBV	1103955	1744211	7.77%	4.453%
10	10.055	1465	1471	1484	rVB	1174780	1556805	6.94%	3.975%
11	11.085	1634	1640	1650	rBV	955540	1238915	5.52%	3.163%
12	12.024	1788	1794	1806	rBV	1358255	1628375	7.26%	4.157%

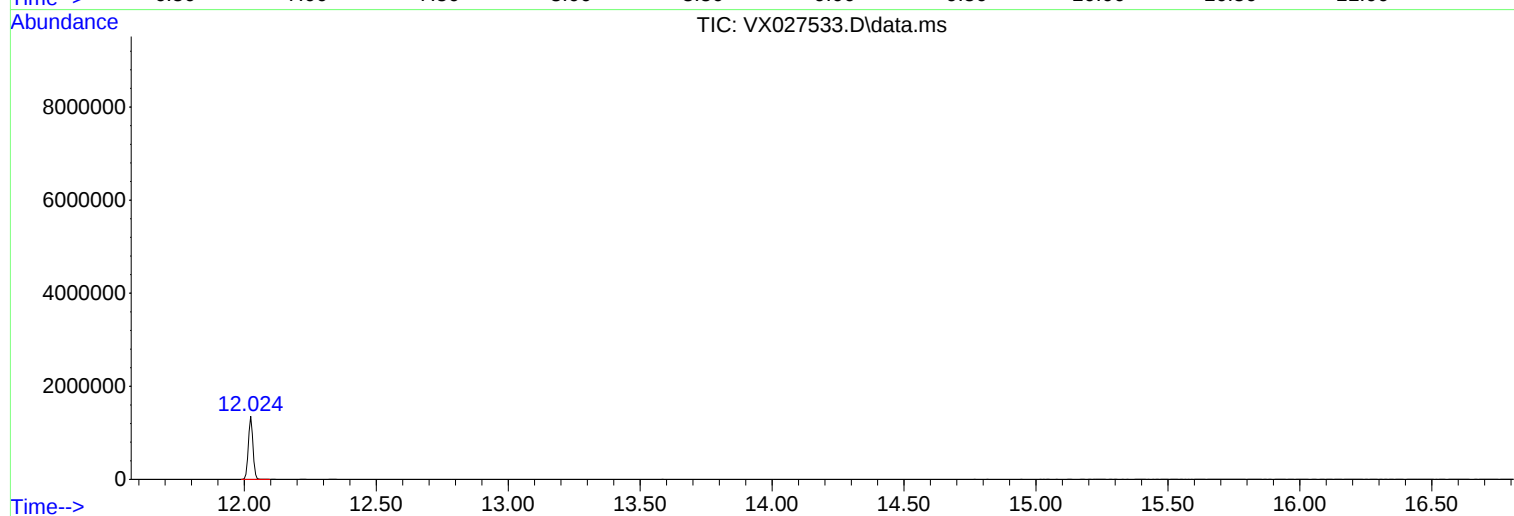
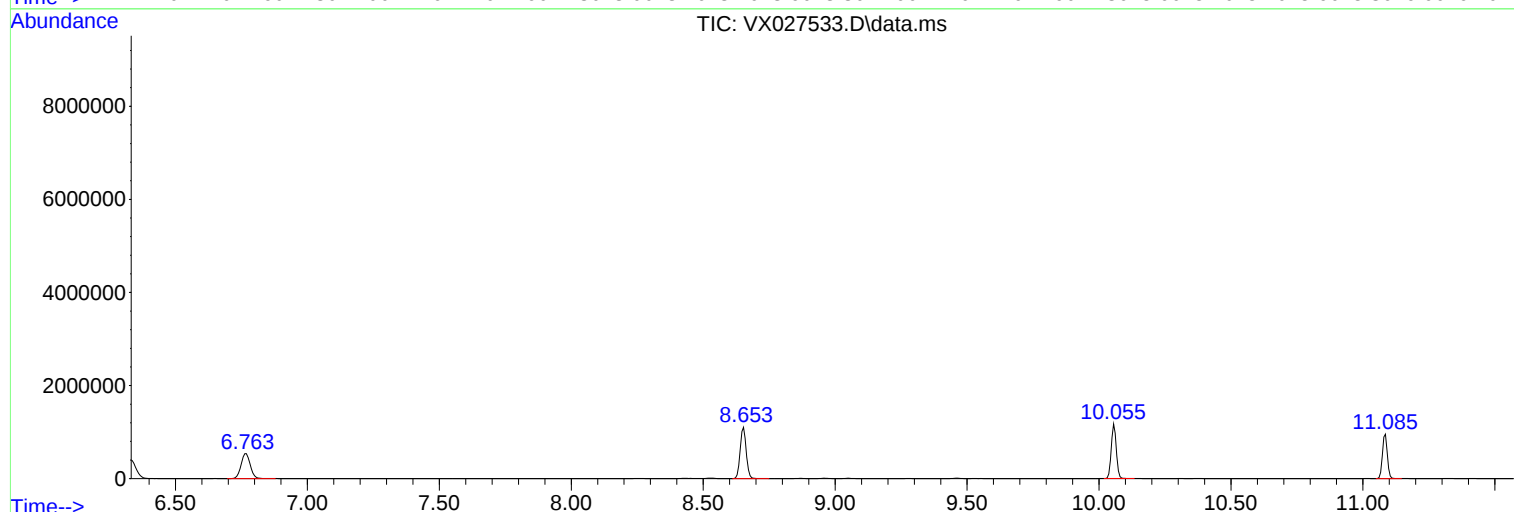
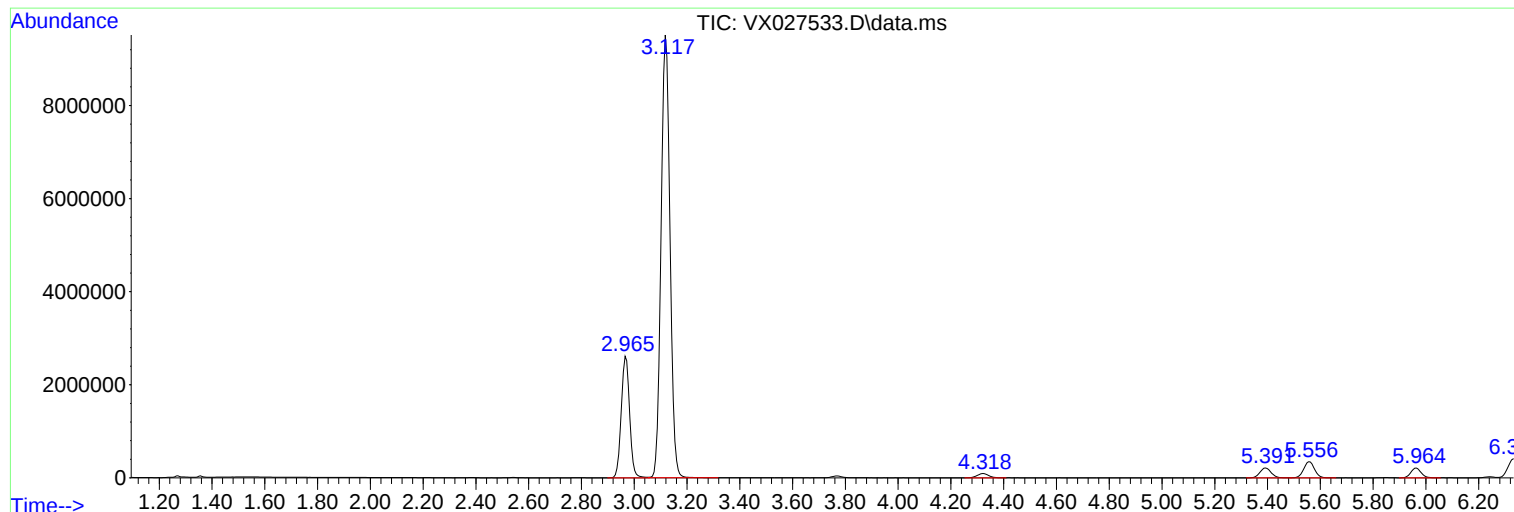
Sum of corrected areas: 39169315

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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X031722W.M
Quant Title : SW846 8260

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



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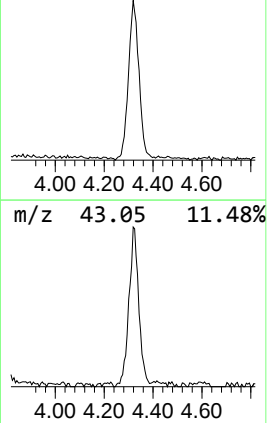
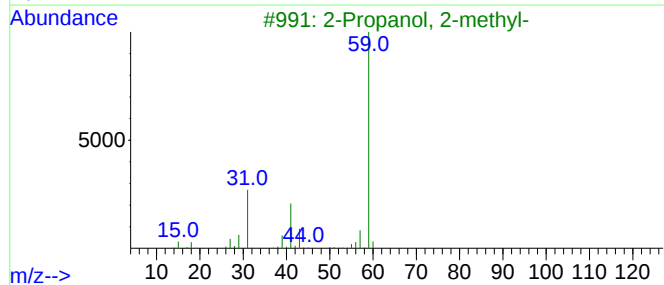
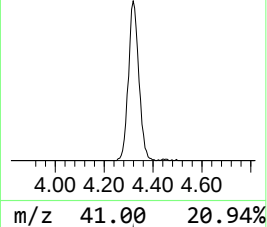
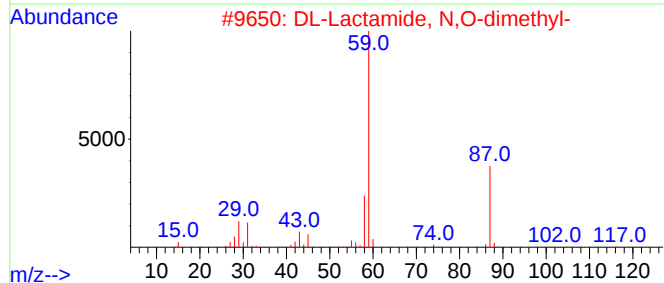
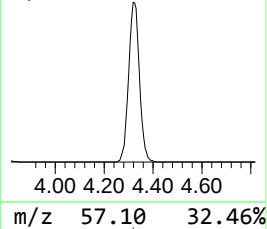
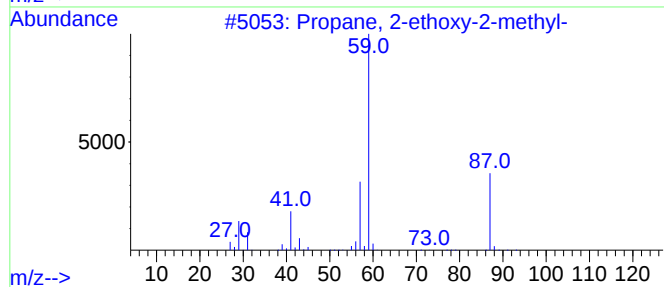
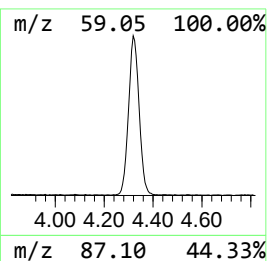
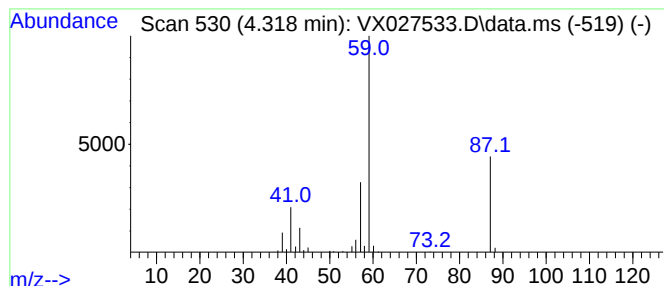
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Peak Number 1 Propane, 2-ethoxy-2-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.318	13.94 ug/l	275445	Pentafluorobenzene	5.562

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propane, 2-ethoxy-2-methyl-	102	C6H14O	000637-92-3	83
2			DL-Lactamide, N,O-dimethyl-	117	C5H11NO2	1000452-56-3	56
3			2-Propanol, 2-methyl-	74	C4H10O	000075-65-0	40
4			2-Pentanol, 2-methyl-	102	C6H14O	000590-36-3	38
5			4-Hydroxy-3-hexanone	116	C6H12O2	004984-85-4	28



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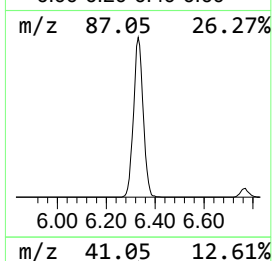
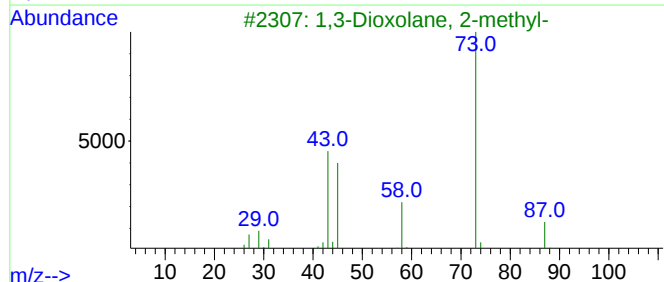
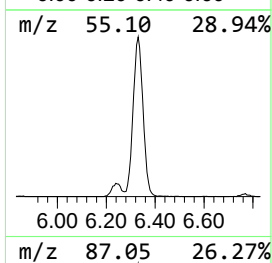
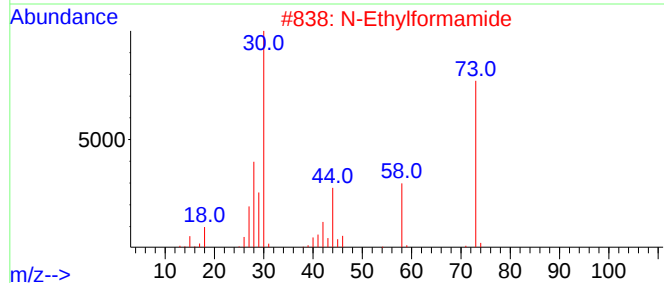
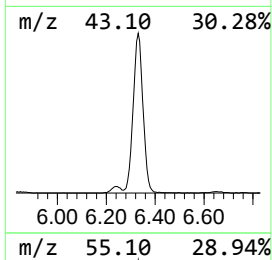
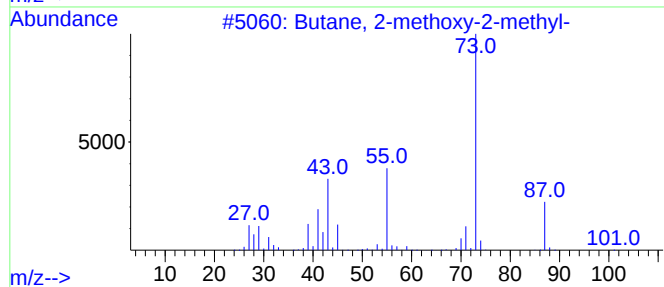
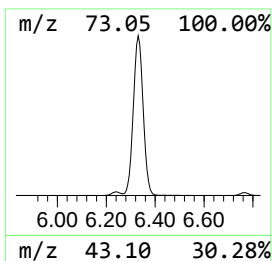
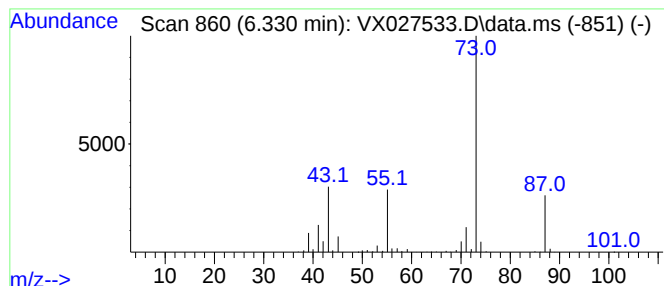
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Peak Number 2 Butane, 2-methoxy-2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.330	43.85 ug/l	1131090	1,4-Difluorobenzene	6.769

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	74
2			N-Ethylformamide	73	C3H7NO	000627-45-2	9
3			1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	9
4			1,4-Butanediamine	88	C4H12N2	000110-60-1	9
5			Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Propane, 2-etho...	4.318	13.9	ug/l	275445	1	5.562	988185	50.0
Butane, 2-metho...	6.330	43.9	ug/l	1131090	2	6.769	1289810	50.0