

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX031822\
 Data File : VX027536.D
 Acq On : 18 Mar 2022 18:28
 Operator : JC/MD
 Sample : N1998-07
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-133

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: VX027536.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.971	298	309	323	rBV	2695874	5894554	32.46%	17.422%
2	3.117	323	333	352	rVB	7807020	18161191	100.00%	53.677%
3	5.391	694	706	721	rBV	213216	616093	3.39%	1.821%
4	5.556	721	733	749	rVB	345117	976350	5.38%	2.886%
5	5.964	789	800	816	rBV	207729	560003	3.08%	1.655%
6	6.239	833	845	852	rBV2	91999	257258	1.42%	0.760%
7	6.330	852	860	875	rVB	149818	426584	2.35%	1.261%
8	6.769	921	932	947	rBV	530534	1271816	7.00%	3.759%
9	8.653	1233	1241	1257	rBV	1075829	1687066	9.29%	4.986%
10	10.055	1465	1471	1485	rVB	1113155	1471499	8.10%	4.349%
11	11.085	1634	1640	1650	rBV	875339	1126043	6.20%	3.328%
12	12.024	1789	1794	1809	rVB	1141184	1385706	7.63%	4.096%

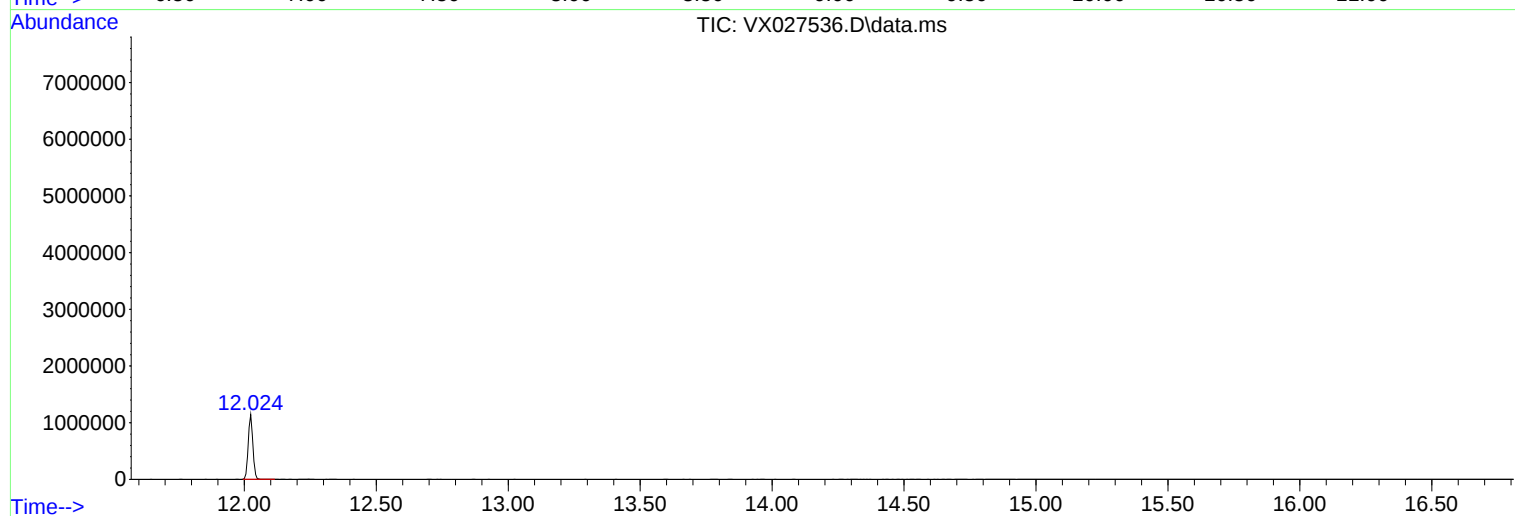
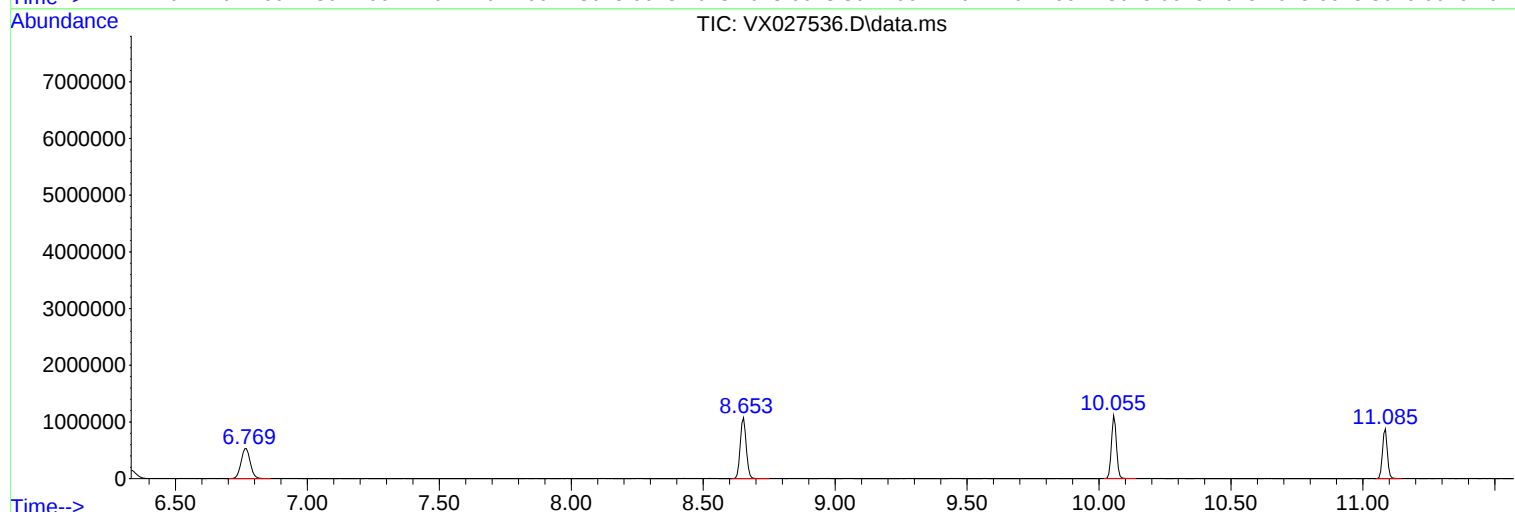
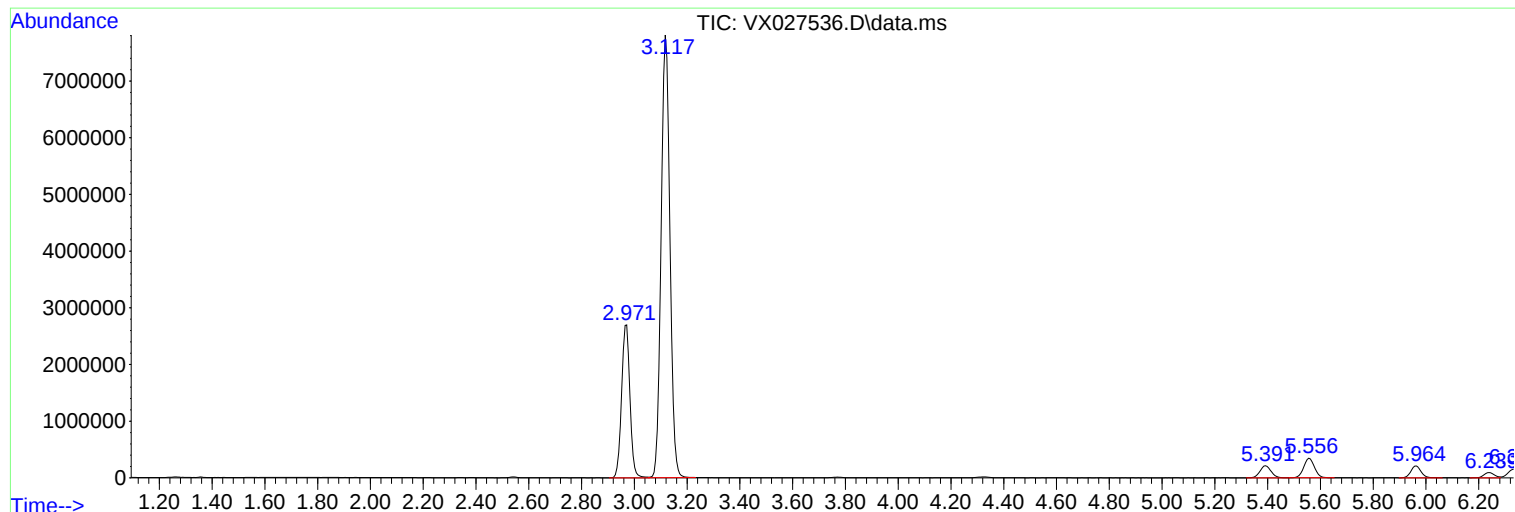
Sum of corrected areas: 33834163

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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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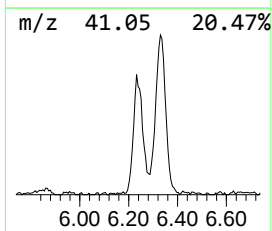
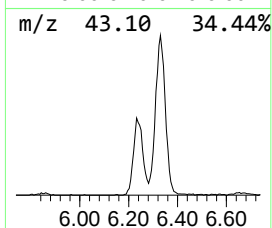
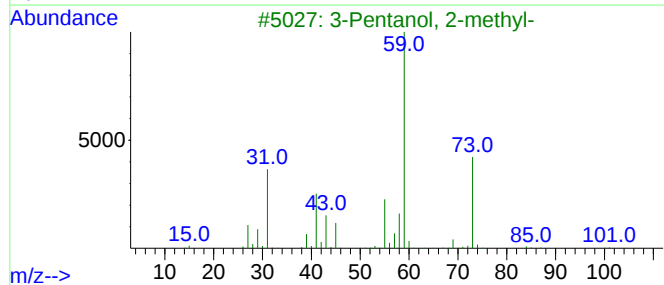
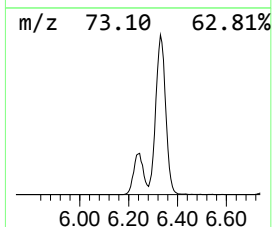
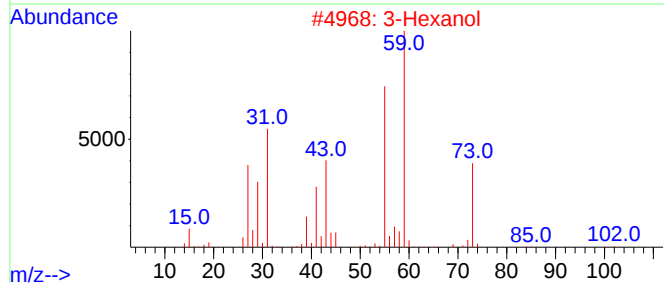
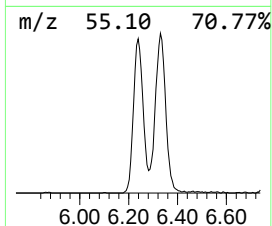
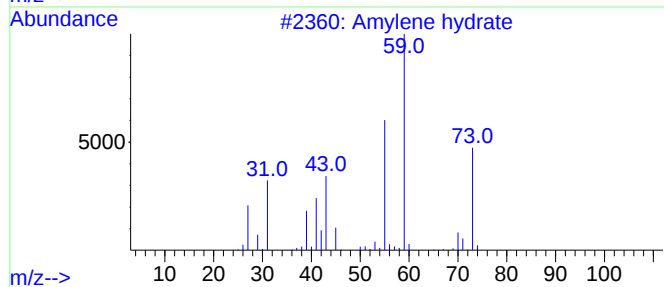
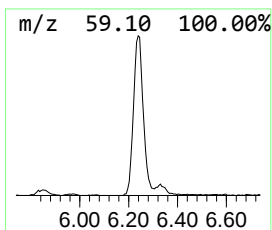
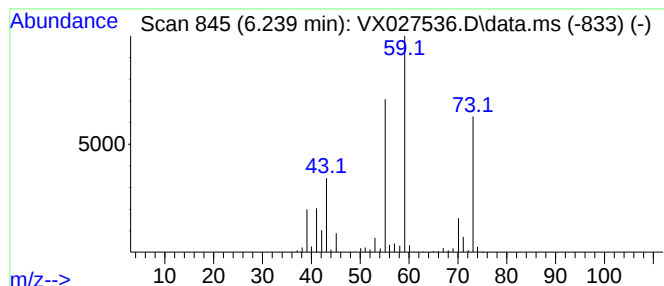
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TIC Integration Parameters: LSCINT.P

Peak Number 1 Amylene hydrate Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.239	10.11 ug/l	257258	1,4-Difluorobenzene	6.769

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Amylene hydrate	88	C5H12O	000075-85-4	74
2			3-Hexanol	102	C6H14O	000623-37-0	40
3			3-Pentanol, 2-methyl-	102	C6H14O	000565-67-3	38
4			Butane, 2-methoxy-	88	C5H12O	006795-87-5	36
5			1-Butene, 3-methyl-	70	C5H10	000563-45-1	11



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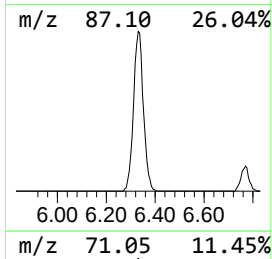
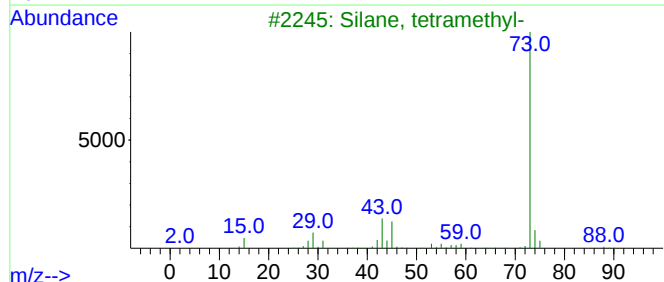
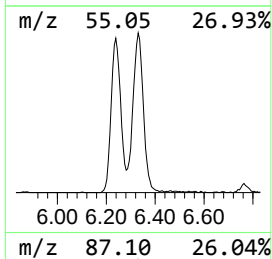
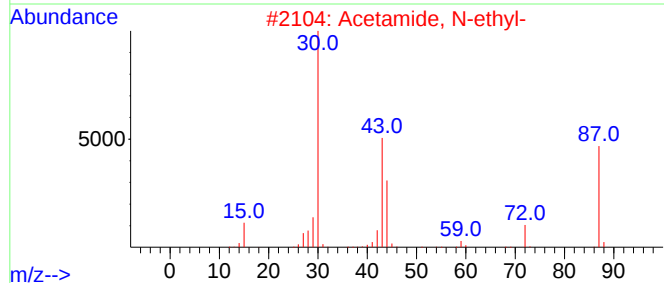
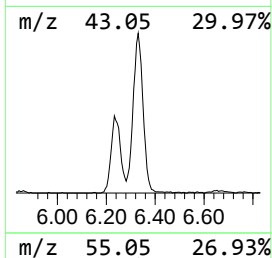
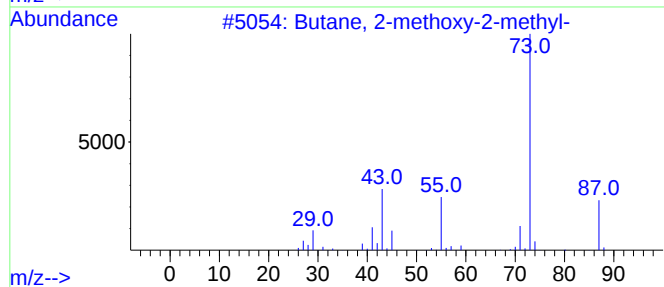
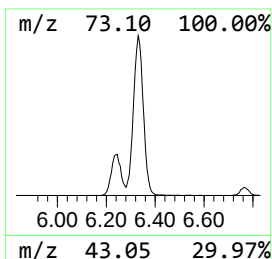
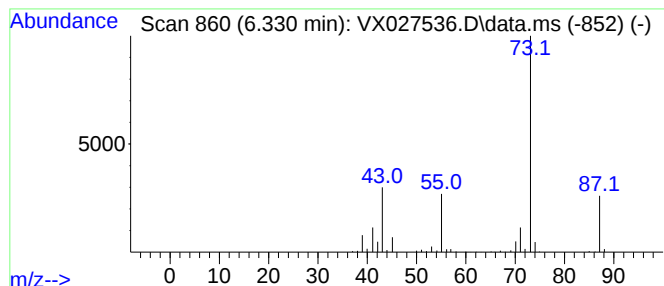
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Peak Number 2 Butane, 2-methoxy-2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.330	16.77 ug/l	426584	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	72
2			Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	9
3			Silane, tetramethyl-	88	C4H12Si	000075-76-3	9
4			1,3-Dioxolane, 2-propyl-	116	C6H12O2	003390-13-4	9
5			1-Propene, 3-methoxy-	72	C4H8O	000627-40-7	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Amylene hydrate	6.239	10.1	ug/l	257258	2	6.769	1271820	50.0
Butane, 2-metho...	6.330	16.8	ug/l	426584	2	6.769	1271820	50.0