

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX121918\
 Data File : VX006644.D
 Acq On : 19 Dec 2018 18:31
 Operator : JC/MD
 Sample : J6479-01
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-1

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_X\METHOD\82X121818W.M

Title : SW846 8260

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.410	36	42	46	rBV2	193822	234460	7.01%	0.932%
2	1.452	46	49	55	rVB	30695	42034	1.26%	0.167%
3	1.605	67	74	80	rBV2	29372	51376	1.54%	0.204%
4	1.696	85	89	92	rVV2	71743	110382	3.30%	0.439%
5	1.733	92	95	103	rVB	434531	574945	17.19%	2.286%
6	1.965	129	133	139	rVB	54320	76536	2.29%	0.304%
7	2.190	164	170	178	rVB	103419	184505	5.51%	0.734%
8	2.379	195	201	212	rBV	2105923	3345526	100.00%	13.303%
9	2.861	276	280	287	rVB	22131	41117	1.23%	0.163%
10	3.702	410	418	432	rVB	303656	738405	22.07%	2.936%
11	4.598	553	565	578	rBV	174991	496872	14.85%	1.976%
12	5.501	703	713	727	rBV	353377	1021745	30.54%	4.063%
13	5.665	729	740	753	rVV	619174	1762963	52.70%	7.010%
14	5.793	754	761	770	rVB6	13223	37979	1.14%	0.151%
15	6.067	796	806	815	rBV2	381727	980892	29.32%	3.900%
16	6.196	822	827	836	rVB	47547	118679	3.55%	0.472%
17	6.372	845	856	868	rBV	416523	1110637	33.20%	4.416%
18	6.860	926	936	955	rBV	941104	2207593	65.99%	8.778%
19	7.214	988	994	997	rBV3	20122	40113	1.20%	0.160%
20	7.262	997	1002	1009	rVB	36857	78415	2.34%	0.312%
21	7.744	1075	1081	1094	rBV	393606	761776	22.77%	3.029%
22	8.043	1123	1130	1139	rBV	32824	62763	1.88%	0.250%
23	8.506	1199	1206	1213	rVB2	42702	79590	2.38%	0.316%
24	8.713	1233	1240	1252	rBV	1954161	3050238	91.17%	12.129%
25	9.512	1364	1371	1377	rBV2	53493	88124	2.63%	0.350%
26	9.792	1412	1417	1421	rVB2	30716	44126	1.32%	0.175%
27	10.109	1464	1469	1479	rVV	1889999	2619251	78.29%	10.415%
28	10.695	1561	1565	1572	rVB	23258	34867	1.04%	0.139%
29	10.902	1595	1599	1604	rBV4	28358	39764	1.19%	0.158%
30	11.134	1632	1637	1649	rBV2	1654804	2271085	67.88%	9.031%
31	11.664	1720	1724	1730	rVB	26447	38574	1.15%	0.153%
32	12.079	1786	1792	1800	rBV	2059048	2543588	76.03%	10.114%
33	12.298	1821	1828	1834	rBV	47075	64799	1.94%	0.258%
34	12.902	1919	1927	1932	rVB3	23508	38258	1.14%	0.152%

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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

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Title : SW846 8260

35	13.347	1995	2000	2005	rBV	118816	156061	4.66%	0.621%
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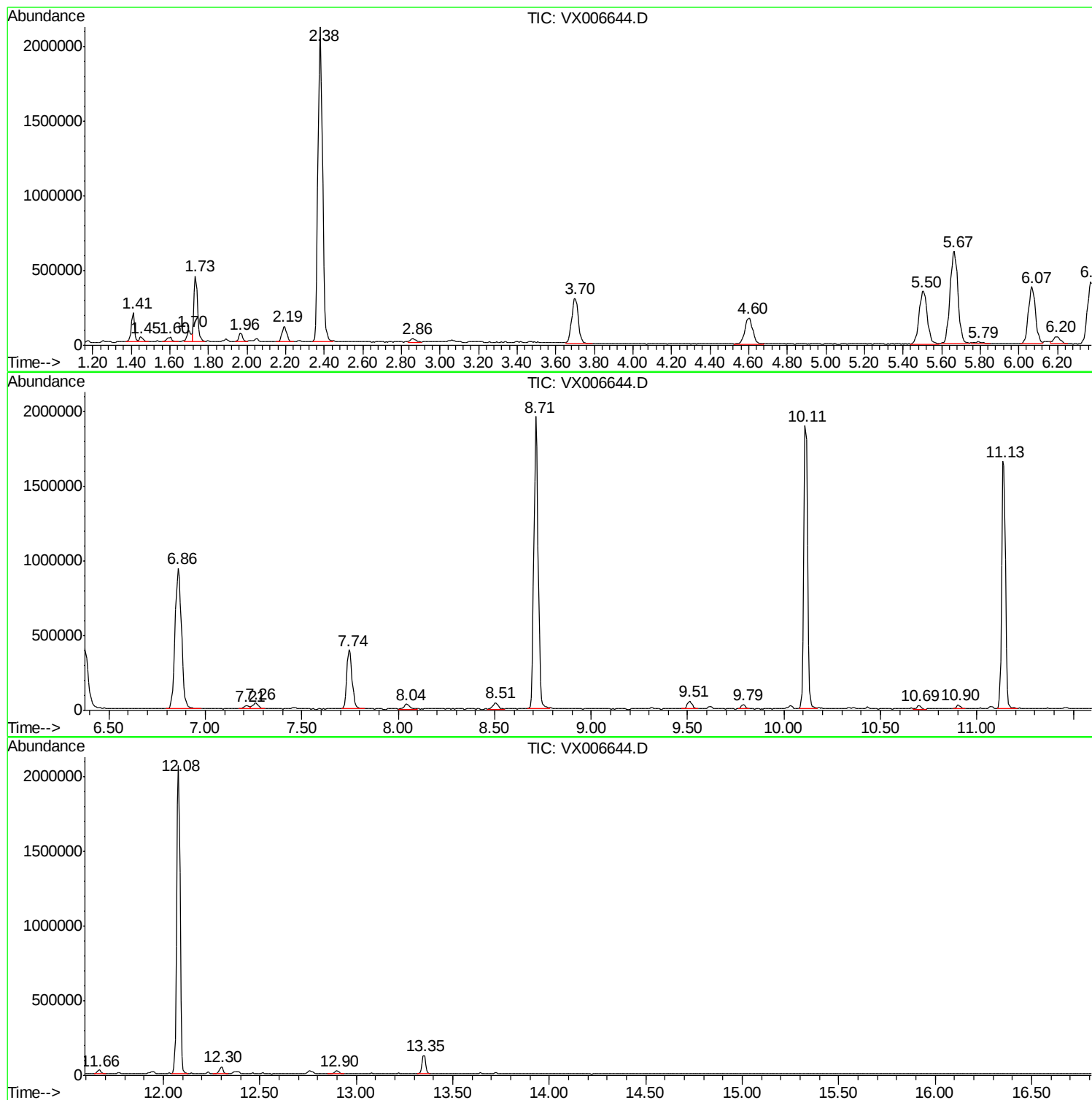
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X121818W.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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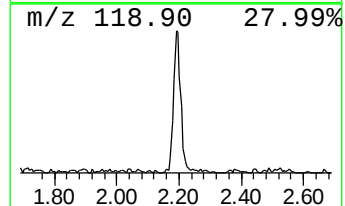
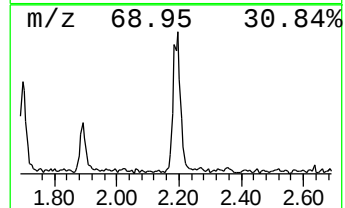
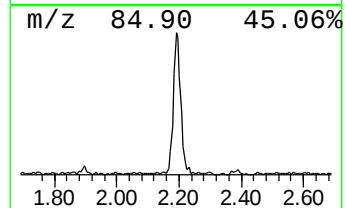
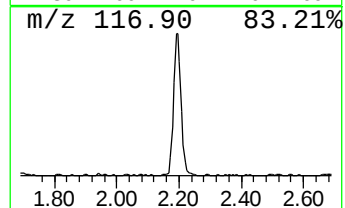
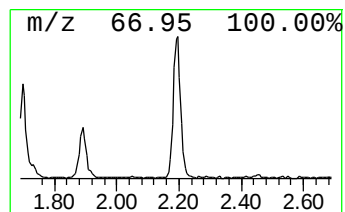
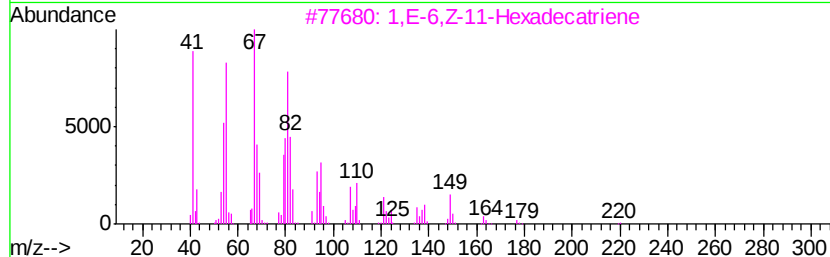
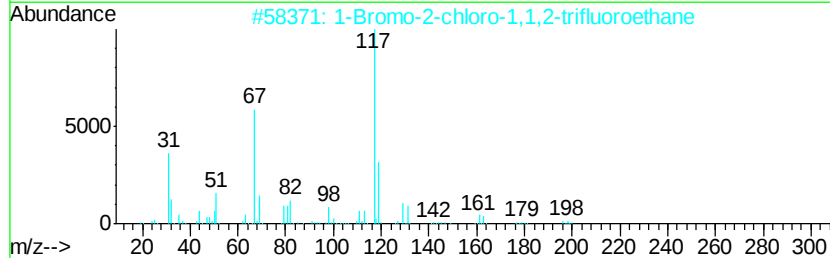
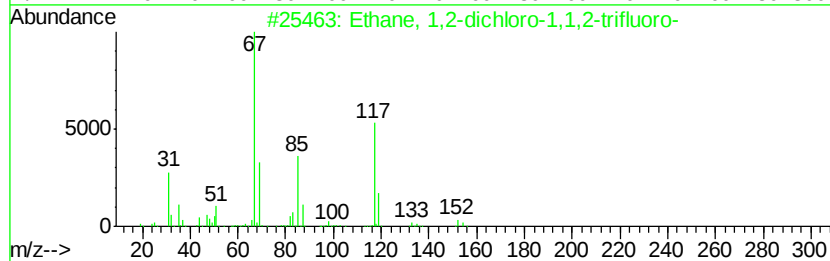
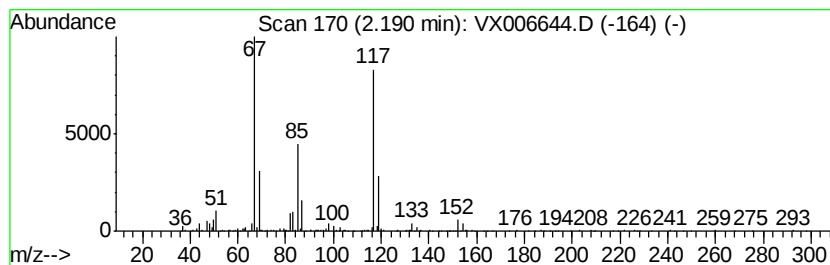
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TIC Integration Parameters: LSCINT.P

Peak Number 1 Ethane, 1,2-dichloro-1,1,2-... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.19	5.23 ug/l	184505	Pentafluorobenzene	5.67

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Ethane, 1,2-dichloro-1,1,2-trifl...	152	C2HCl2F3	000354-23-4	94
2		1-Bromo-2-chloro-1,1,2-trifluoro...	196	C2HBrClF3	000354-06-3	45
3		1,E-6,Z-11-Hexadecatriene	220	C16H28	083483-55-0	18
4		Phosphorochloridic acid, hexyl p...	270	C11H24ClO3P	1000309-08-9	12
5		1,4-Pentadiene, 3-methyl-	82	C6H10	001115-08-8	10



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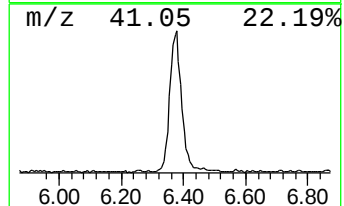
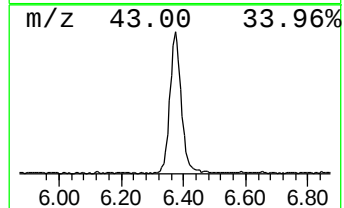
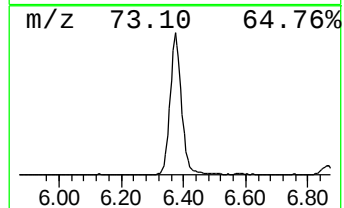
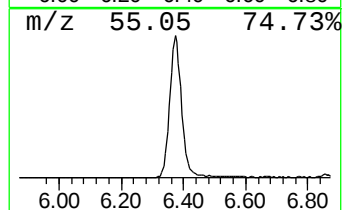
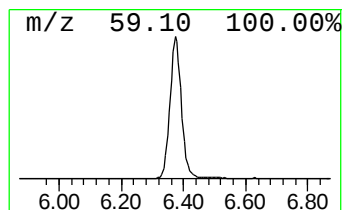
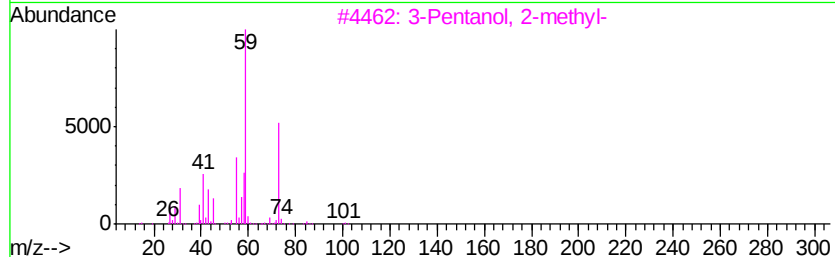
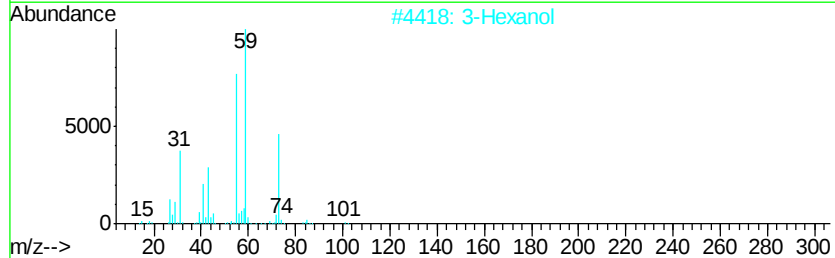
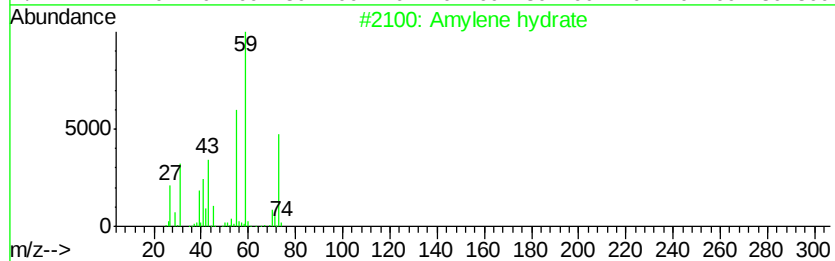
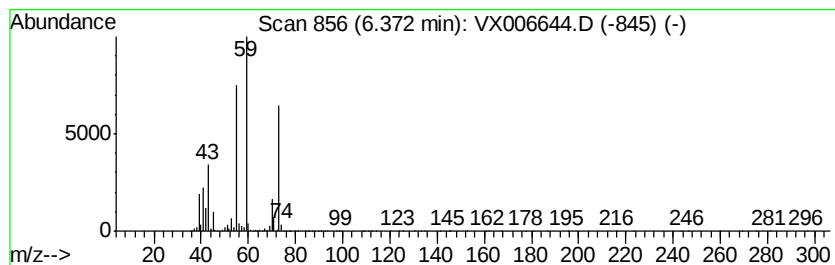
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Peak Number 2 Amylene hydrate Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.37	25.15 ug/l	1110640	1,4-Difluorobenzene	6.86

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Amylene hydrate	88	C5H12O	000075-85-4	90
2		3-Hexanol	102	C6H14O	000623-37-0	45
3		3-Pentanol, 2-methyl-	102	C6H14O	000565-67-3	42
4		Silane, trimethyl-	74	C3H10Si	000993-07-7	32
5		Cyclopropane, 1,2-dimethyl-, cis-	70	C5H10	000930-18-7	18



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Ethane, 1,2-dichl...	2.19	5.2	ug/l	184505	1	5.67	1762960	50.0
Amylene hydrate	6.37	25.1	ug/l	1110640	2	6.86	2207590	50.0