

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX062119\
 Data File : VX010383.D
 Acq On : 21 Jun 2019 14:38
 Operator : JC/SP
 Sample : K3473-01
 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\82X061919W.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	1.282	18	21	26	rVB	11318	16945	1.14%	0.145%
2	1.410	36	42	46	rBV	95755	114173	7.66%	0.979%
3	1.446	46	48	53	rVB	14990	18101	1.22%	0.155%
4	1.696	85	89	91	rBV2	35157	43796	2.94%	0.376%
5	1.733	91	95	103	rVB	148958	206241	13.85%	1.769%
6	1.964	127	133	138	rBV	26672	38866	2.61%	0.333%
7	2.190	164	170	178	rVB	49699	83750	5.62%	0.718%
8	2.269	178	183	191	rVB3	7890	18064	1.21%	0.155%
9	2.373	193	200	210	rBV	903300	1489587	100.00%	12.777%
10	2.611	233	239	249	rVB	17368	33926	2.28%	0.291%
11	3.043	298	310	318	rVB3	7356	22602	1.52%	0.194%
12	3.702	407	418	430	rBV	139132	341464	22.92%	2.929%
13	4.598	555	565	576	rBV	93274	256992	17.25%	2.204%
14	5.500	701	713	725	rBV	162126	469299	31.51%	4.025%
15	5.665	729	740	750	rVV	277052	791512	53.14%	6.789%
16	5.750	752	754	767	rVB2	8340	18803	1.26%	0.161%
17	6.061	796	805	815	rBV	154546	403404	27.08%	3.460%
18	6.195	821	827	835	rVB2	20600	53929	3.62%	0.463%
19	6.342	840	851	866	rBV	269579	720425	48.36%	6.180%
20	6.860	925	936	951	rBV	419051	975158	65.46%	8.365%
21	7.220	986	995	997	rBV3	11088	24483	1.64%	0.210%
22	7.256	997	1001	1008	rVB2	18713	40310	2.71%	0.346%
23	7.732	1072	1079	1092	rBV	220681	436569	29.31%	3.745%
24	8.043	1124	1130	1137	rBV2	18590	34565	2.32%	0.296%
25	8.494	1197	1204	1211	rBV	29914	51918	3.49%	0.445%
26	8.713	1233	1240	1249	rBV	879214	1343836	90.22%	11.527%
27	9.506	1362	1370	1378	rBV2	35558	62480	4.19%	0.536%
28	9.792	1410	1417	1424	rBV	18745	27698	1.86%	0.238%
29	10.030	1447	1456	1460	rBV3	14308	25558	1.72%	0.219%
30	10.109	1464	1469	1479	rVV	814721	1133137	76.07%	9.720%
31	10.902	1594	1599	1604	rBV2	21222	30927	2.08%	0.265%
32	11.073	1622	1627	1631	rVB2	10126	15445	1.04%	0.132%
33	11.140	1631	1638	1647	rBV2	745091	1040509	69.85%	8.925%
34	11.664	1720	1724	1729	rBV	18733	25946	1.74%	0.223%

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35	12.078	1786	1792	1800	rBV	828239	1069309	71.79%	9.172%
36	12.298	1825	1828	1834	rVB	34263	43813	2.94%	0.376%
37	12.755	1899	1903	1911	rVB5	11357	21581	1.45%	0.185%
38	12.901	1919	1927	1932	rBV2	12486	22031	1.48%	0.189%
39	13.346	1995	2000	2006	rBV	69404	91060	6.11%	0.781%

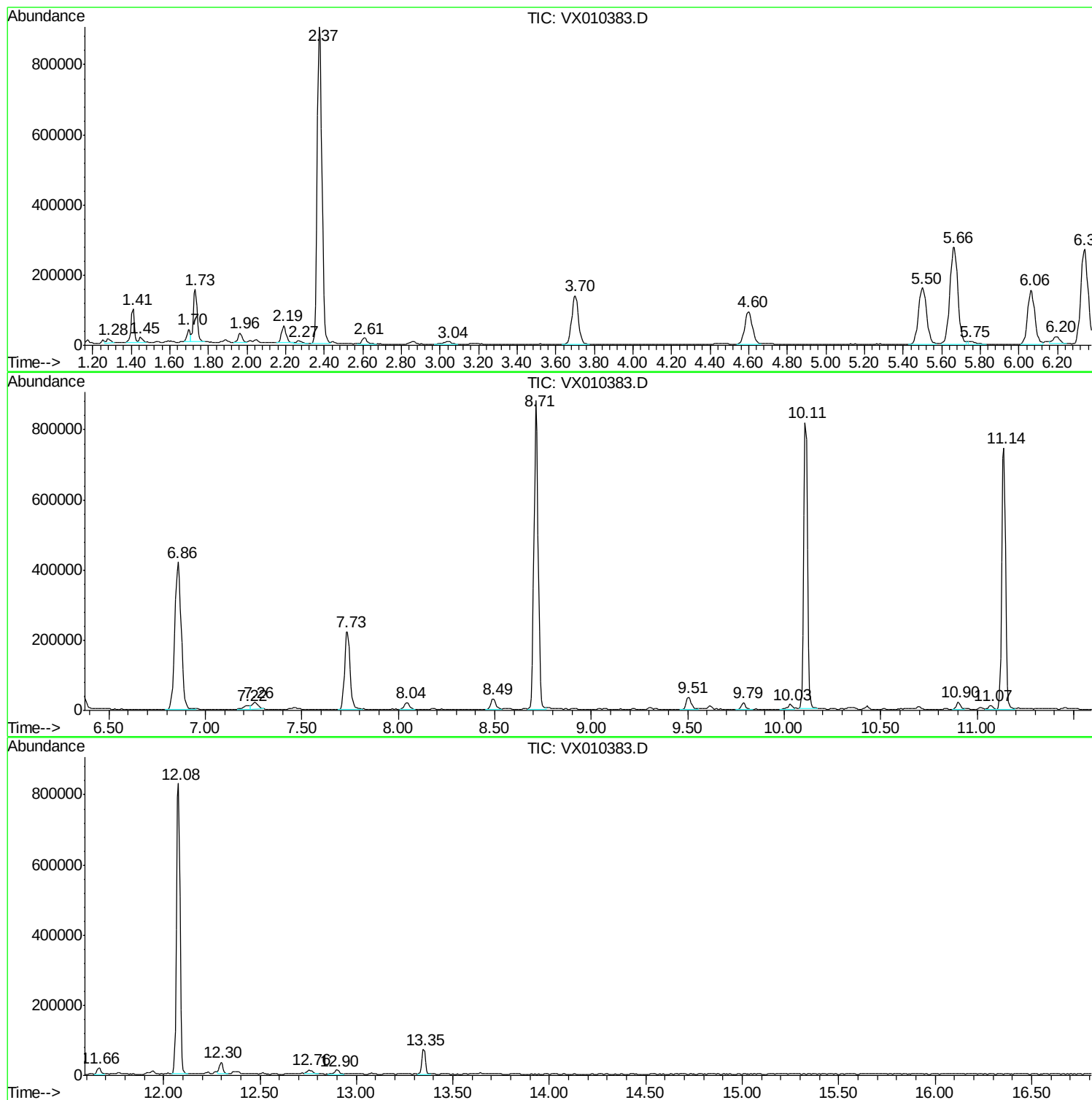
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
Quant Title : SW846 8260

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



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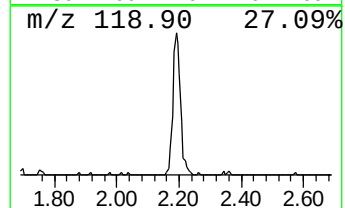
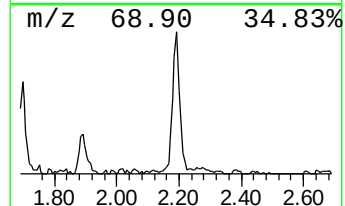
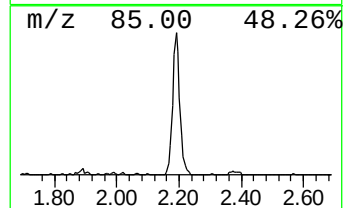
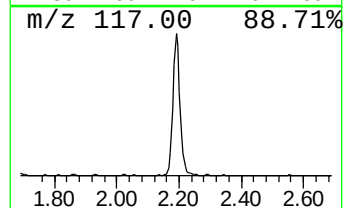
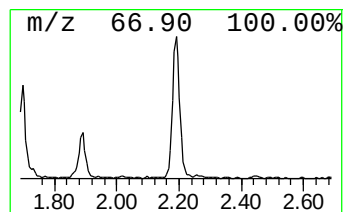
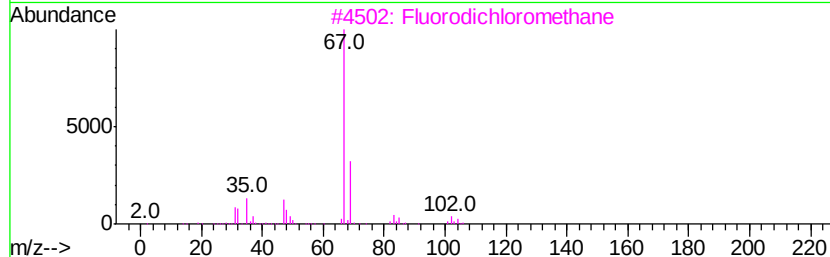
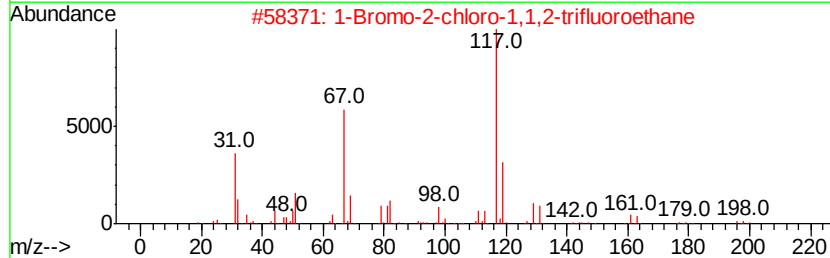
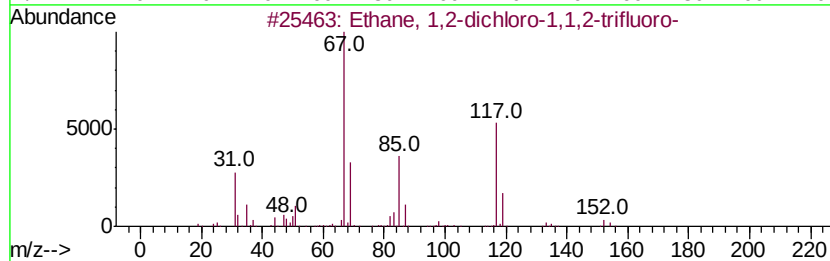
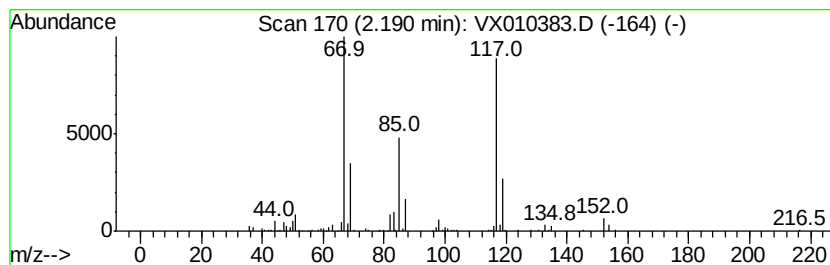
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Peak Number 1 Ethane, 1,2-dichloro-1,1,2-... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.19	5.29 ug/l	83750	Pentafluorobenzene	5.66

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	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	47
3		Fluorodichloromethane	102	CHCl2F	000075-43-4	16
4		Cyclopropane, 1,2-dimethyl-3-met...	82	C6H10	062338-02-7	10
5		Methyl 2-butynoate	98	C5H6O2	023326-27-4	9



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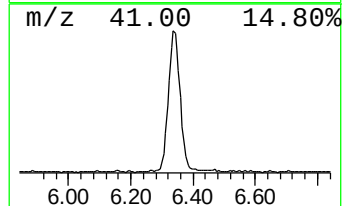
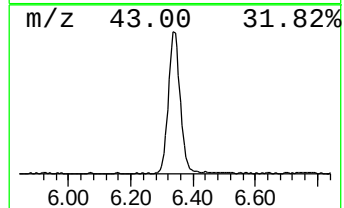
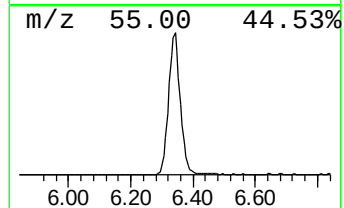
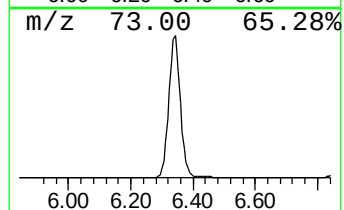
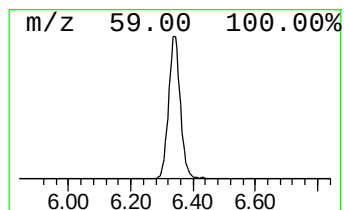
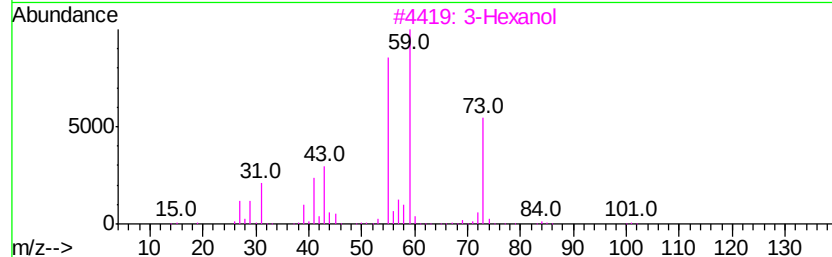
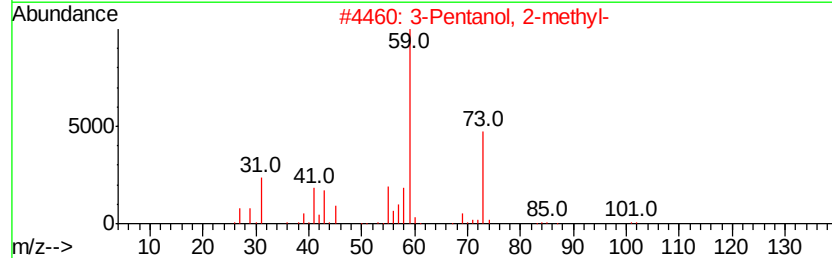
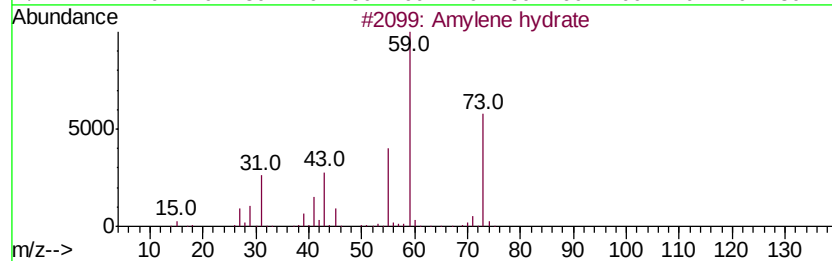
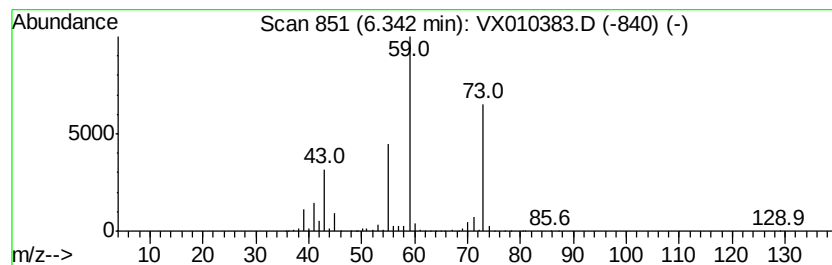
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Peak Number 2 Amylene hydrate Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.34	36.94 ug/l	720425	1,4-Difluorobenzene	6.86

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Amylene hydrate	88	C5H12O	000075-85-4	83
2		3-Pentanol, 2-methyl-	102	C6H14O	000565-67-3	40
3		3-Hexanol	102	C6H14O	000623-37-0	39
4		Propanoic acid, 2-hydroxy-2-methyl-	104	C4H8O3	000594-61-6	33
5		2-Methyl-5-hexen-3-ol	114	C7H14O	032815-70-6	10



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Ethane, 1,2-dichl...	2.19	5.3	ug/l	83750	1	5.66	791512	50.0
Amylene hydrate	6.34	36.9	ug/l	720425	2	6.86	975158	50.0