

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX080320\
 Data File : VX017753.D
 Acq On : 03 Aug 2020 18:26
 Operator : JC/SP
 Sample : L3536-01
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 JTY1-GW-01

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\82X072320W.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	1.160	9	12	21	rVB	365707	321721	11.14%	1.430%
2	1.380	44	48	53	rBV2	103832	130283	4.51%	0.579%
3	1.477	60	64	70	rVB	156725	173462	6.01%	0.771%
4	2.331	200	204	208	rBV	17305	29575	1.02%	0.131%
5	2.416	212	218	226	rVV	153877	248023	8.59%	1.102%
6	2.568	237	243	252	rVB2	84606	170773	5.91%	0.759%
7	2.837	281	287	295	rVB	149530	270007	9.35%	1.200%
8	4.306	517	528	538	rBV2	41618	119471	4.14%	0.531%
9	4.647	577	584	593	rBV5	13318	44361	1.54%	0.197%
10	5.470	708	719	731	rBV2	355739	1017120	35.23%	4.520%
11	5.635	733	746	763	rVB	646933	1816196	62.90%	8.072%
12	6.037	800	812	818	rBV	386226	1027909	35.60%	4.568%
13	6.117	818	825	841	rVB	1101650	2887395	100.00%	12.833%
14	6.836	933	943	953	rBV	889218	2123241	73.53%	9.436%
15	7.641	1069	1075	1082	rVB3	16255	34903	1.21%	0.155%
16	8.695	1240	1248	1255	rBV	1849628	2850754	98.73%	12.670%
17	8.769	1255	1260	1266	rVB	273723	423935	14.68%	1.884%
18	9.561	1384	1390	1397	rBV2	88657	127817	4.43%	0.568%
19	10.098	1470	1478	1493	rBV	1898314	2520425	87.29%	11.202%
20	10.238	1495	1501	1507	rVB	40871	56800	1.97%	0.252%
21	10.341	1514	1518	1523	rVB	48458	67635	2.34%	0.301%
22	10.689	1569	1575	1585	rVB7	22124	56582	1.96%	0.251%
23	10.896	1604	1609	1613	rBV	43551	55120	1.91%	0.245%
24	11.122	1641	1646	1655	rBV	1639307	1996445	69.14%	8.873%
25	11.457	1696	1701	1705	rVB2	33968	45697	1.58%	0.203%
26	11.963	1781	1784	1787	rBV4	21307	30722	1.06%	0.137%
27	12.000	1787	1790	1794	rVB	55692	67883	2.35%	0.302%
28	12.061	1794	1800	1808	rVB	2071274	2547060	88.21%	11.320%
29	12.158	1811	1816	1825	rVB	71612	100101	3.47%	0.445%
30	12.256	1827	1832	1847	rBV	622055	964775	33.41%	4.288%
31	12.823	1913	1925	1930	rBV	16833	49016	1.70%	0.218%
32	12.890	1930	1936	1941	rVV3	57002	86662	3.00%	0.385%
33	13.036	1950	1960	1964	rBV	19994	38587	1.34%	0.171%

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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

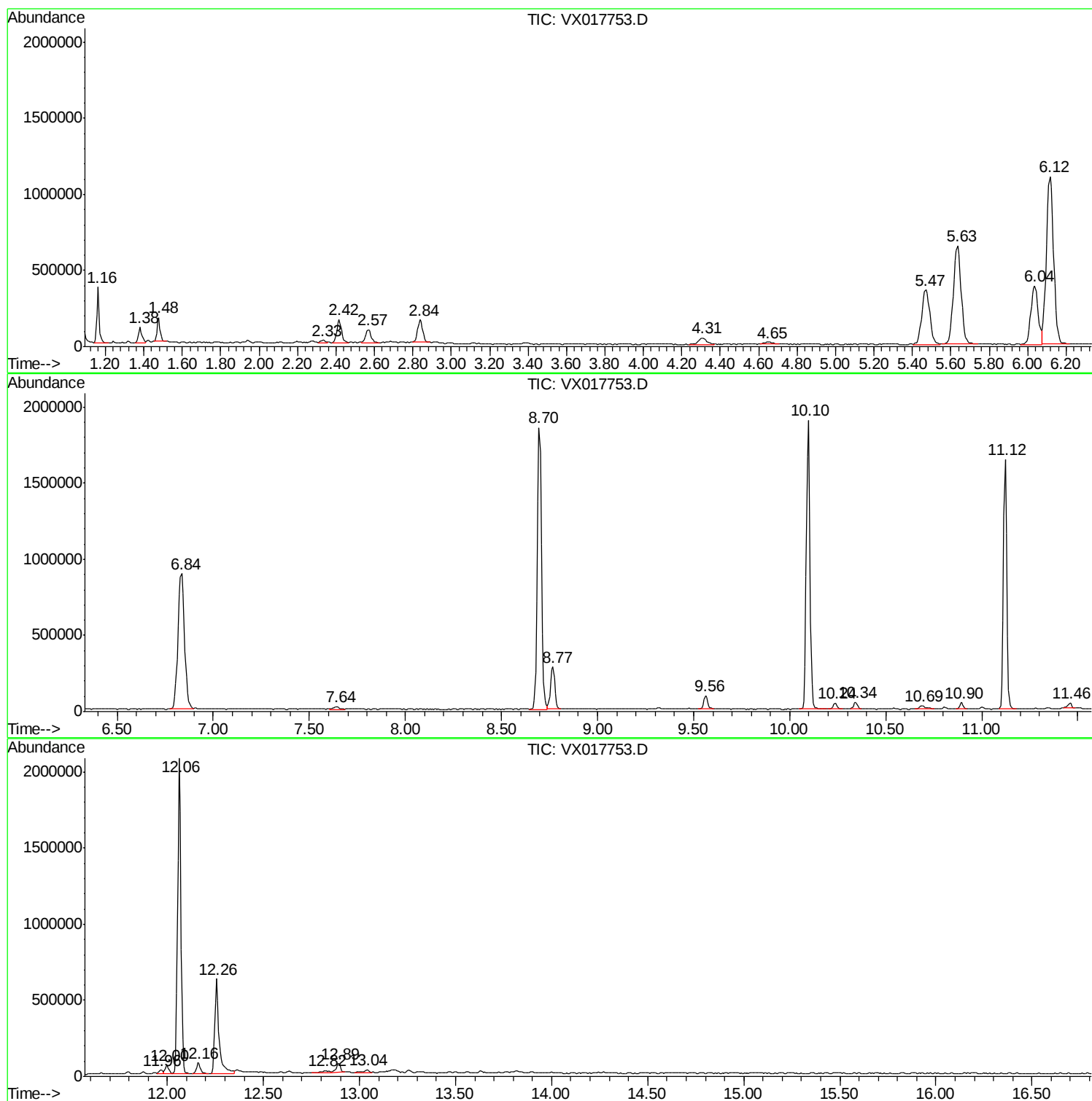
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X072320W.M
Quant Title : SW846 8260

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



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Instrument :
MSVOA_X
ClientSampled :
JTY1-GW-01

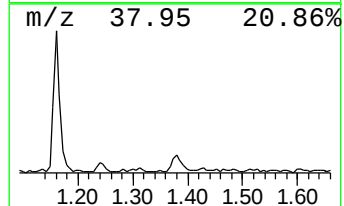
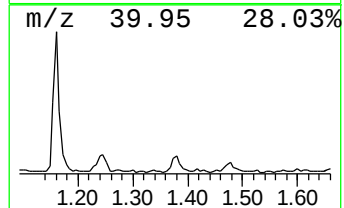
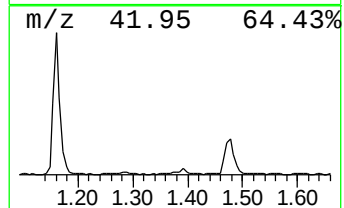
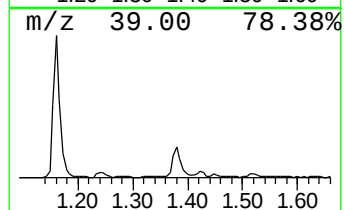
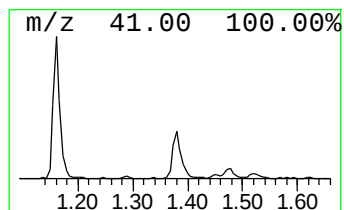
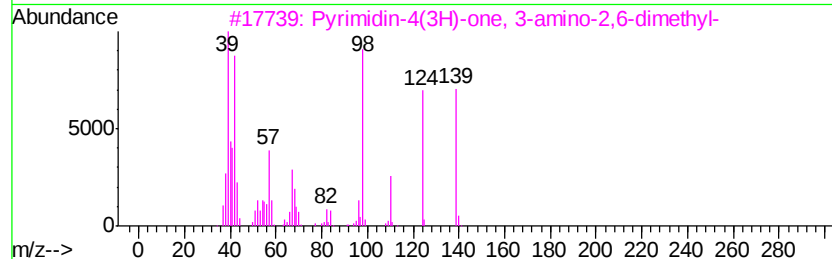
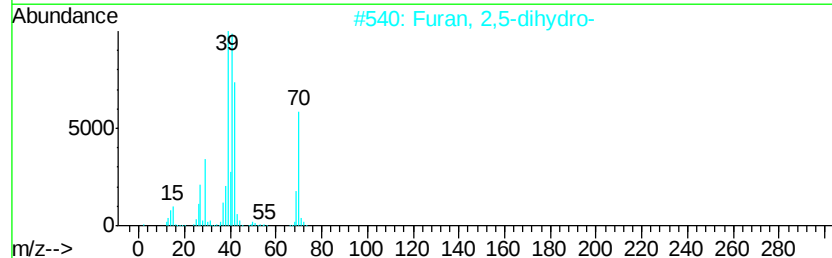
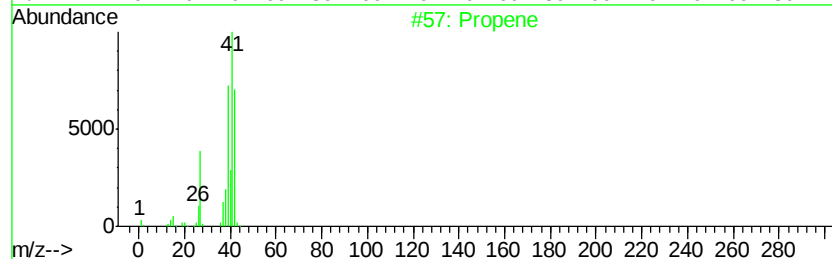
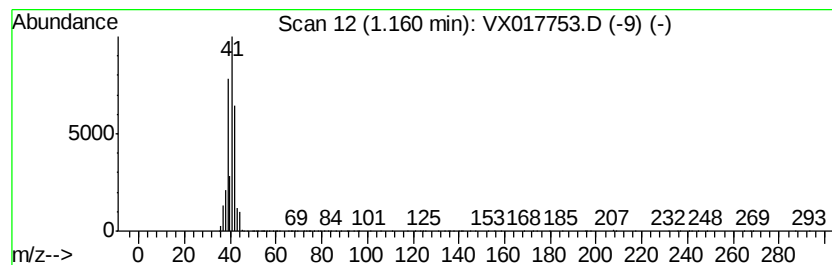
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Peak Number 1 Propene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.16	8.86 ug/l	321721	Pentafluorobenzene	5.63

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Propene	42	C3H6	000115-07-1	90
2		Furan, 2,5-dihydro-	70	C4H6O	001708-29-8	83
3		Pyrimidin-4(3H)-one, 3-amino-2,6...	139	C6H9N3O	093098-74-9	78
4		2-Butenal, (E)-	70	C4H6O	000123-73-9	78
5		Aluminum, tripropyl-	156	C9H21Al	000102-67-0	74



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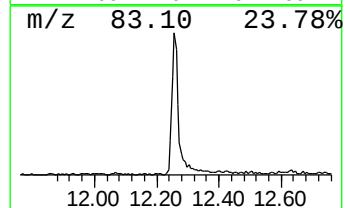
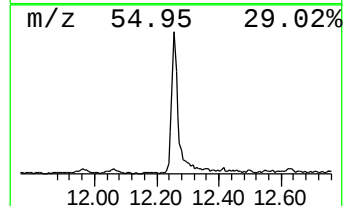
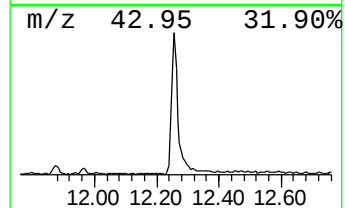
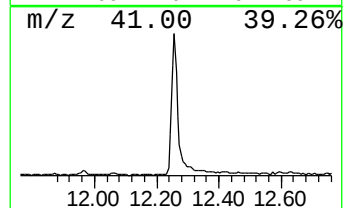
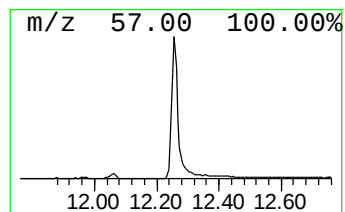
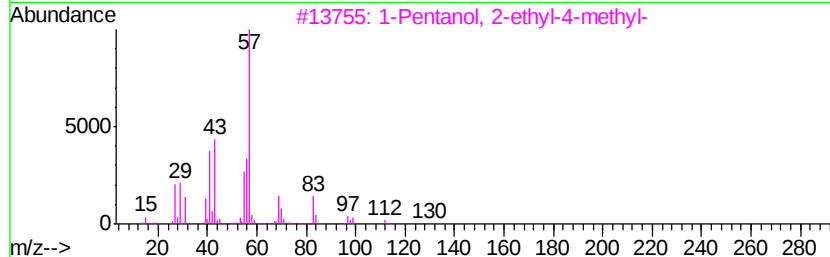
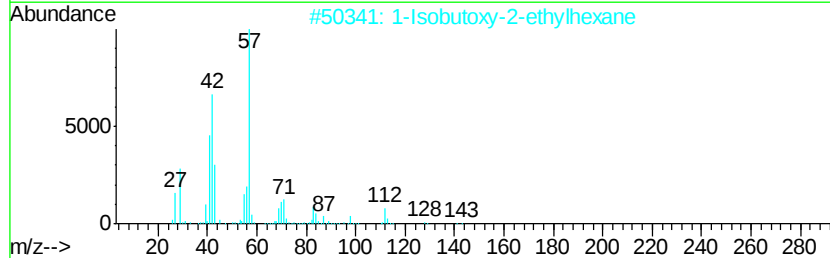
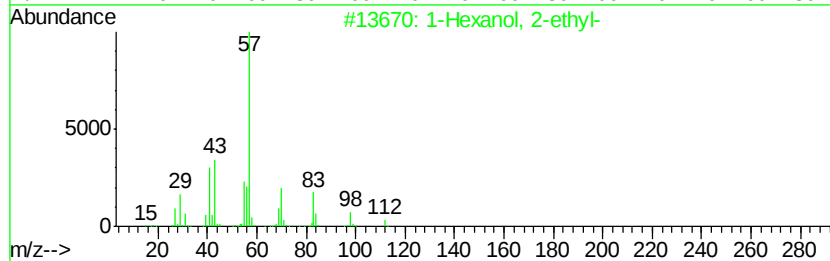
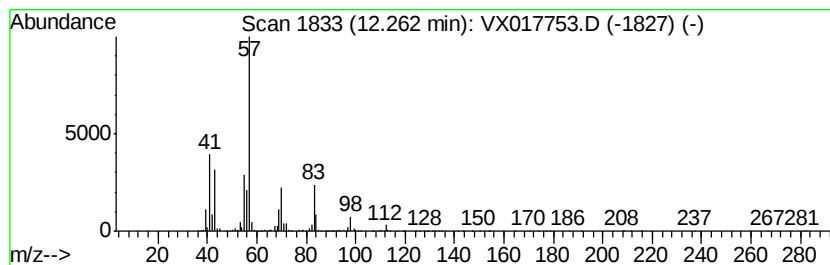
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Peak Number 2 1-Hexanol, 2-ethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.26	18.94 ug/l	964775	1,4-Dichlorobenzene-d4	12.06
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	1-Hexanol, 2-ethyl-	130 C8H18O	000104-76-7	83
2	1-Isobutoxy-2-ethylhexane	186 C12H26O	1000139-90-3	72
3	1-Pentanol, 2-ethyl-4-methyl-	130 C8H18O	000106-67-2	59
4	Oxalic acid, 2-ethylhexyl octyl ...	314 C18H34O4	1000309-39-1	53
5	Heptane, 1,1'-oxybis-	214 C14H30O	000629-64-1	53



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Propene	1.16	8.9	ug/l	321721	1	5.63	1816200	50.0
1-Hexanol, 2-ethyl-	12.26	18.9	ug/l	964775	4	12.06	2547060	50.0