

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX061920\
 Data File : VX016908.D
 Acq On : 20 Jun 2020 05:07
 Operator : JC/SP
 Sample : L3063-07
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 41 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 TW-6B

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\82X061820W.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	142113	133231	8.55%	1.478%
2	1.380	44	48	57	rBV2	67241	99954	6.41%	1.109%
3	1.940	136	140	144	rBV	14607	20712	1.33%	0.230%
4	1.989	144	148	155	rVB3	12076	24887	1.60%	0.276%
5	2.416	213	218	229	rVB	11575	22085	1.42%	0.245%
6	2.574	236	244	252	rBV2	15961	34961	2.24%	0.388%
7	4.647	577	584	585	rBV2	8345	16128	1.03%	0.179%
8	5.464	706	718	732	rBV2	182787	534999	34.32%	5.936%
9	5.629	734	745	761	rVV	291956	824099	52.86%	9.144%
10	6.031	800	811	825	rBV	186015	487736	31.29%	5.412%
11	6.830	932	942	962	rBV	441199	1066869	68.43%	11.838%
12	7.153	987	995	999	rBV2	9836	22573	1.45%	0.250%
13	8.695	1240	1248	1257	rBV	936783	1429941	91.72%	15.867%
14	10.097	1472	1478	1489	rBV	1143765	1558985	100.00%	17.299%
15	11.122	1640	1646	1662	rVB	999704	1273061	81.66%	14.126%
16	12.061	1795	1800	1809	rVB	1143304	1429760	91.71%	15.865%
17	12.256	1827	1832	1838	rBV	20538	32136	2.06%	0.357%

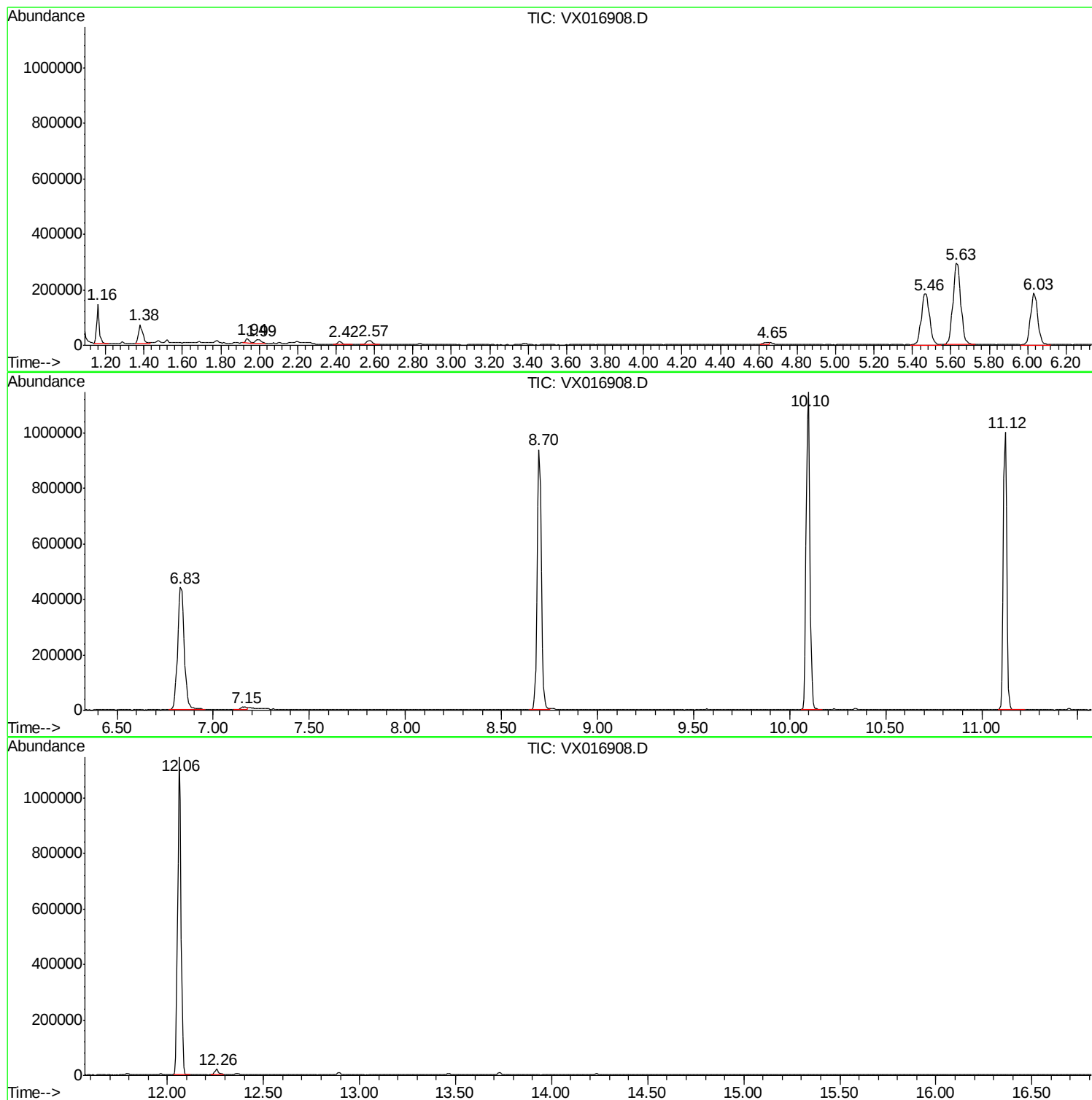
Sum of corrected areas: 9012117

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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X061820W.M
Quant Title : SW846 8260

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



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ClientSampled :
TW-6B

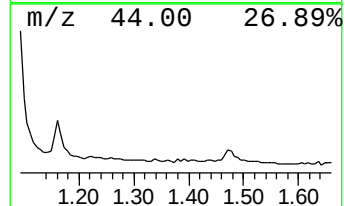
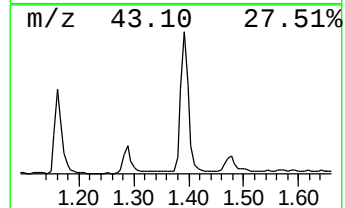
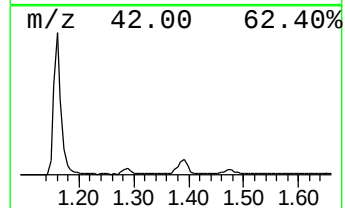
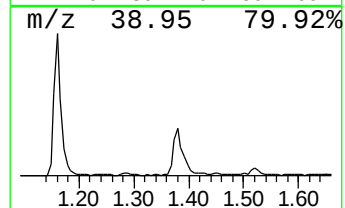
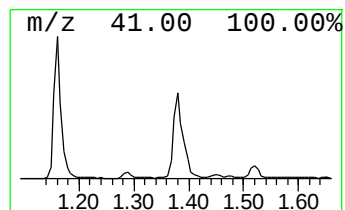
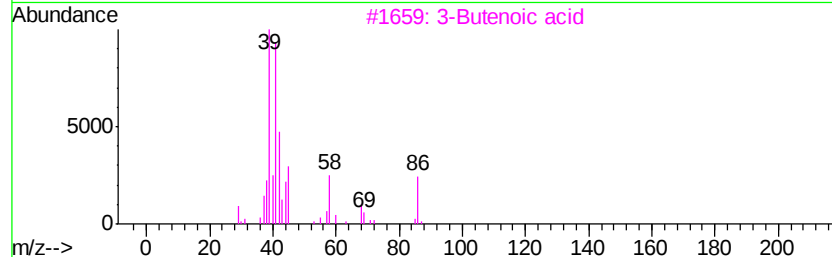
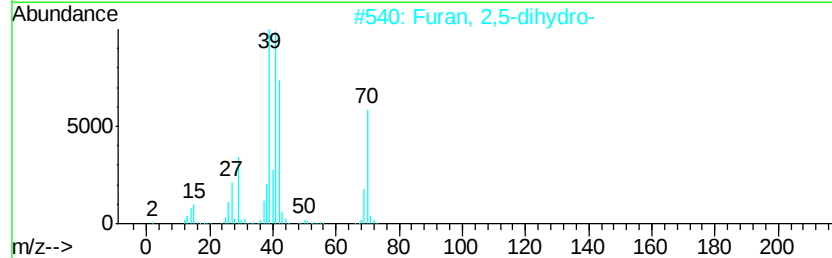
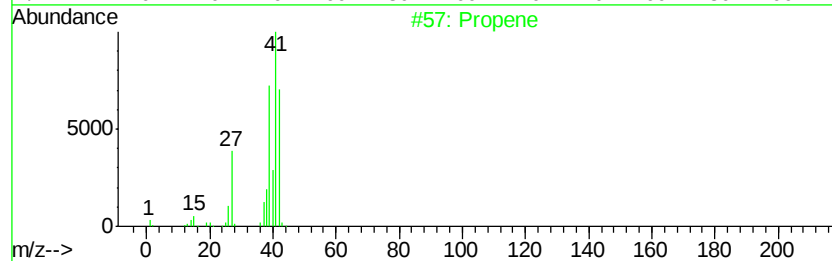
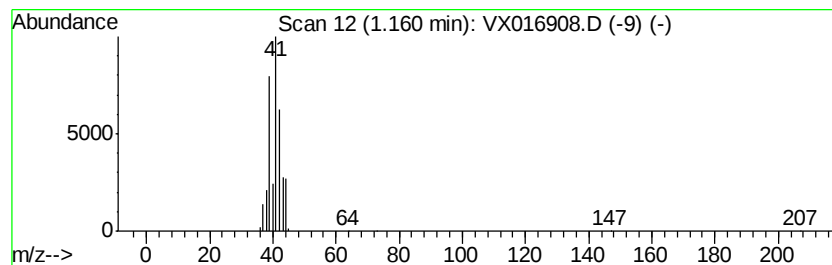
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Peak Number 1 Propene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.16	8.08 ug/l	133231	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		3-Butenoic acid	86	C4H6O2	000625-38-7	83
4		Pyrimidin-4(3H)-one, 3-amino-2,6...	139	C6H9N3O	093098-74-9	78
5		2-Butenal, (E)-	70	C4H6O	000123-73-9	78



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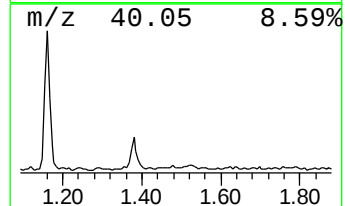
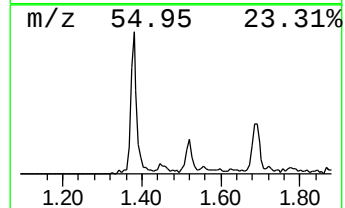
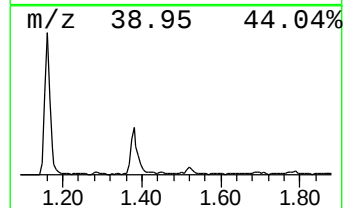
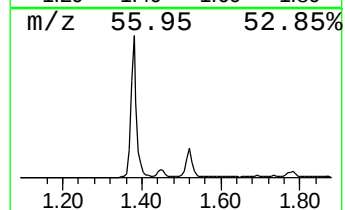
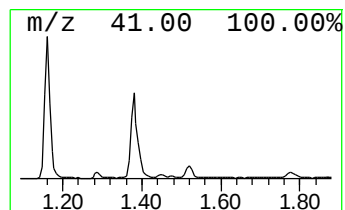
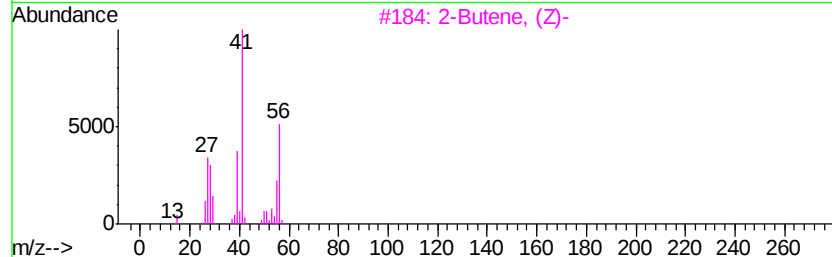
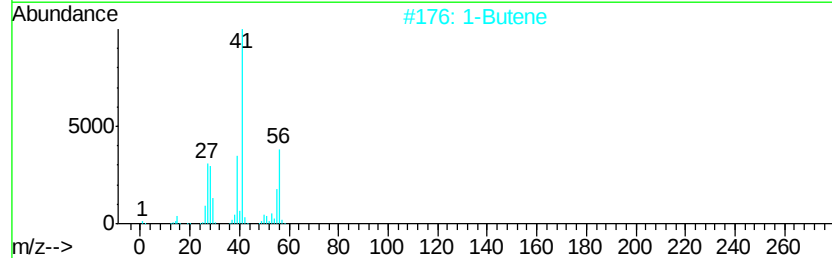
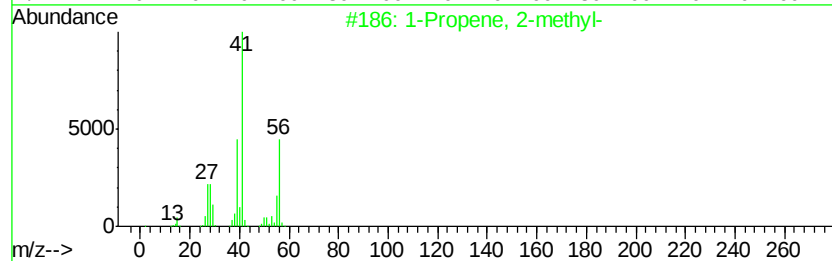
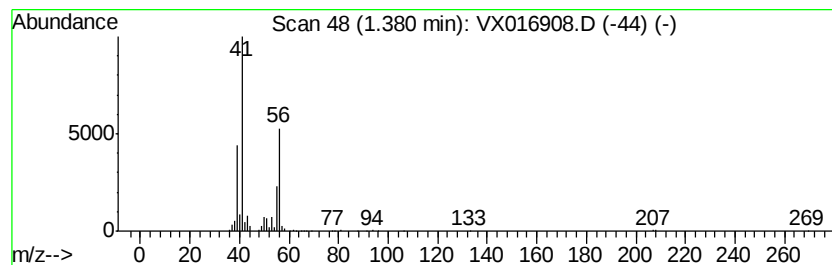
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Peak Number 2 1-Propene, 2-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.38	6.06 ug/l	99954	Pentafluorobenzene	5.63

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Propene, 2-methyl-	56	C4H8	000115-11-7	87
2		1-Butene	56	C4H8	000106-98-9	83
3		2-Butene, (Z)-	56	C4H8	000590-18-1	80
4		2-Butene, (E)-	56	C4H8	000624-64-6	80
5		2-Butene	56	C4H8	000107-01-7	64



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Propene	1.16	8.1	ug/l	133231	1	5.63	824099	50.0
1-Propene, 2-methyl-	1.38	6.1	ug/l	99954	1	5.63	824099	50.0