

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX060320\
 Data File : VX016564.D
 Acq On : 03 Jun 2020 13:41
 Operator : JC/SP
 Sample : L2867-01
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 SP20-GMZ5-ER1

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\82X052720W.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	1.288	29	33	43	rBV	30823	32122	2.34%	0.417%
2	1.374	44	47	58	rVV	137229	134344	9.80%	1.742%
3	1.471	60	63	71	rBV	20154	25302	1.84%	0.328%
4	2.416	213	218	228	rVB	14563	24202	1.76%	0.314%
5	2.575	236	244	253	rBV	25939	48577	3.54%	0.630%
6	5.470	705	719	732	rBV	144347	416181	30.35%	5.396%
7	5.629	734	745	760	rVB	263741	750645	54.74%	9.733%
8	6.031	800	811	828	rBV	159430	426566	31.10%	5.531%
9	6.830	931	942	955	rBV	408485	974299	71.04%	12.633%
10	7.635	1067	1074	1082	rBV2	9934	19771	1.44%	0.256%
11	8.695	1241	1248	1256	rBV	903485	1371415	100.00%	17.782%
12	10.098	1472	1478	1495	rVB	905297	1243785	90.69%	16.127%
13	11.122	1640	1646	1656	rBV	805373	1007772	73.48%	13.067%
14	12.061	1795	1800	1809	rBV	1006571	1237249	90.22%	16.043%

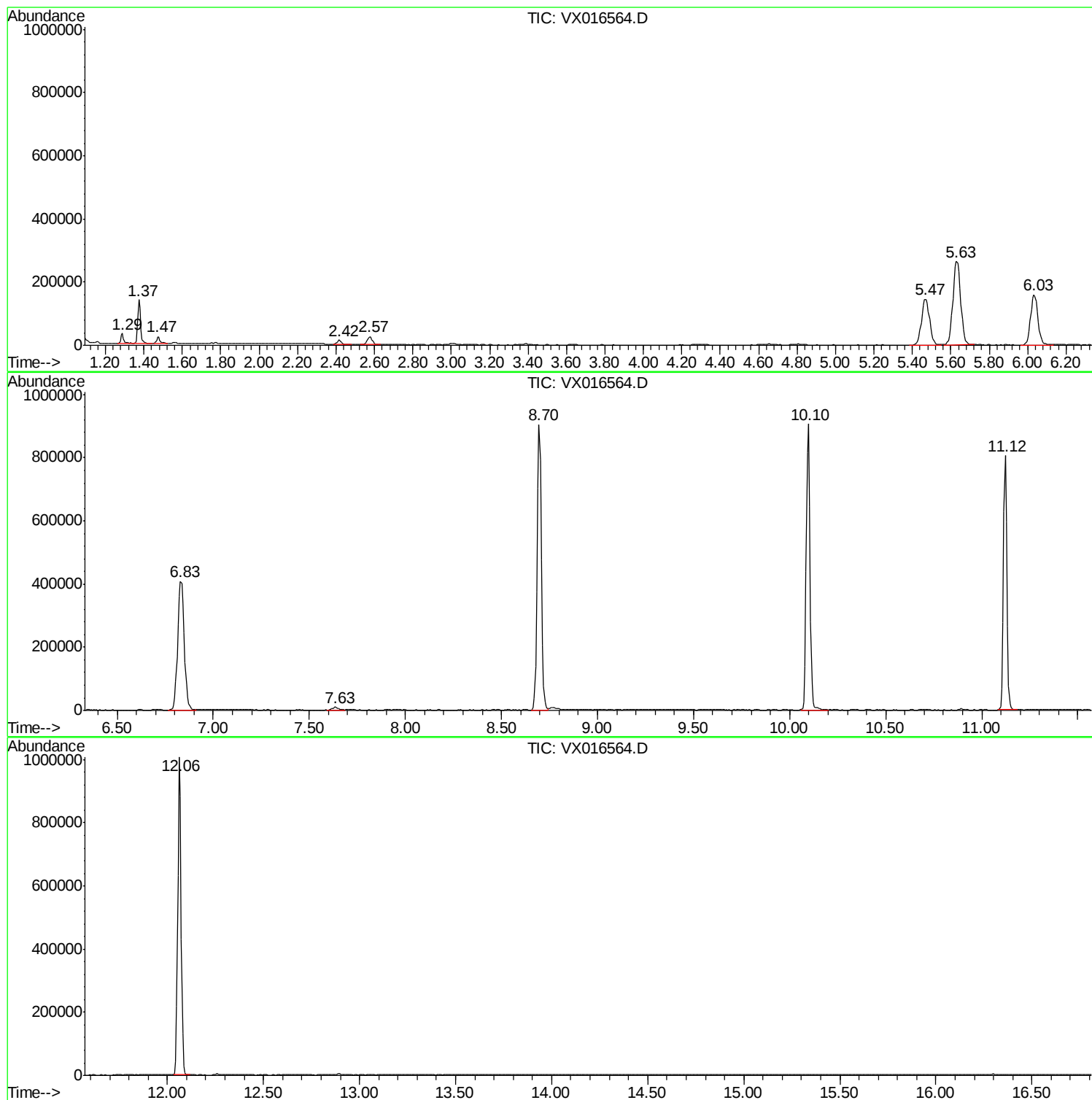
Sum of corrected areas: 7712230

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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X052720W.M
Quant Title : SW846 8260

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



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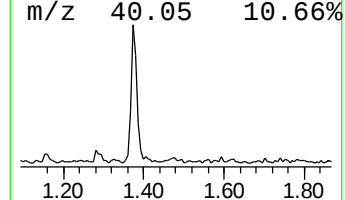
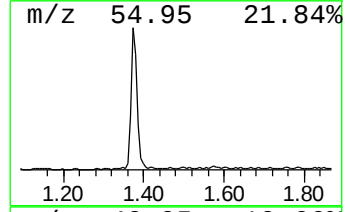
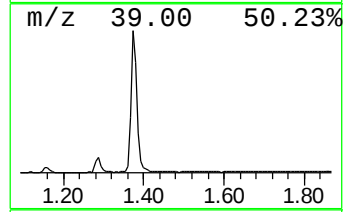
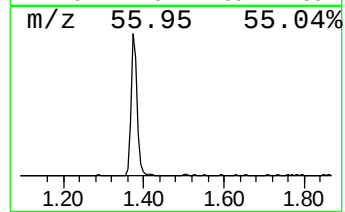
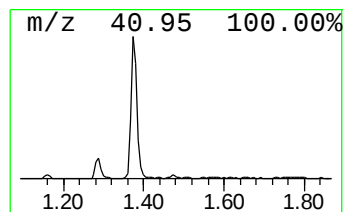
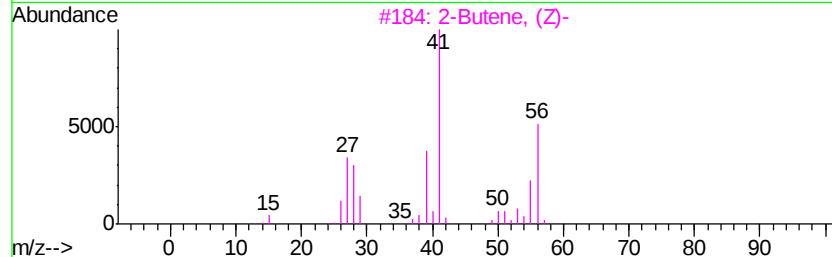
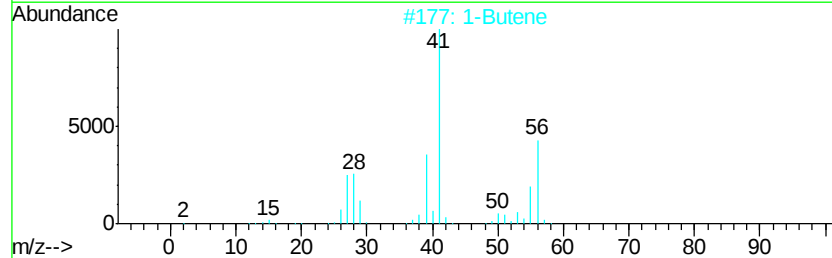
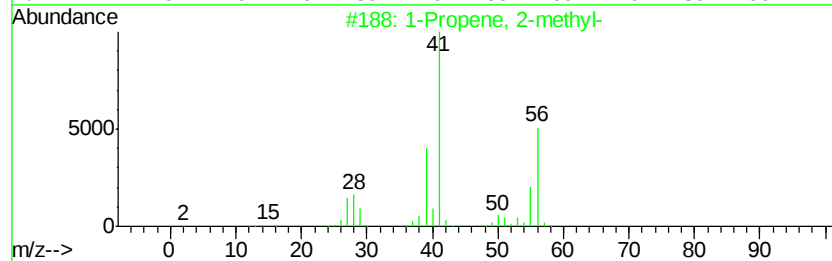
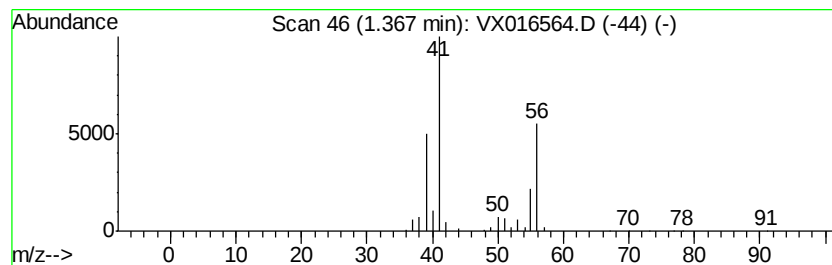
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Peak Number 1 1-Propene, 2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.37	8.95 ug/l	134344	Pentafluorobenzene	5.63

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Propene, 2-methyl-	56	C4H8	000115-11-7	90
2			1-Butene	56	C4H8	000106-98-9	87
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
1-Propene, 2-methyl-	1.37	8.9	ug/l	134344	1	5.63	750645	50.0