

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX100918\
 Data File : VX005194.D
 Acq On : 10 Oct 2018 07:53
 Operator : JC/MD
 Sample : J5303-11
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 42 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 SA-TW514-4044

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\82X092618W.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	1.391	35	39	47	rBV2	36074	57740	5.36%	1.004%
2	1.623	67	77	80	rVV3	9383	21600	2.00%	0.376%
3	1.952	126	131	135	rBV	10079	14010	1.30%	0.244%
4	1.995	135	138	147	rVB4	7473	12797	1.19%	0.223%
5	2.623	235	241	249	rVB	27299	51264	4.76%	0.891%
6	5.500	699	713	728	rBV	139358	399007	37.03%	6.938%
7	5.671	728	741	757	rBV2	175896	514212	47.73%	8.942%
8	6.067	795	806	822	rBV	154846	393725	36.54%	6.846%
9	6.860	925	936	958	rBV	269266	667285	61.93%	11.603%
10	8.713	1232	1240	1261	rBV	640501	1077426	100.00%	18.735%
11	10.115	1464	1470	1488	rBV	603814	840418	78.00%	14.614%
12	11.140	1632	1638	1650	rBV	596847	764251	70.93%	13.289%
13	12.079	1786	1792	1804	rBV	771830	937105	86.98%	16.295%

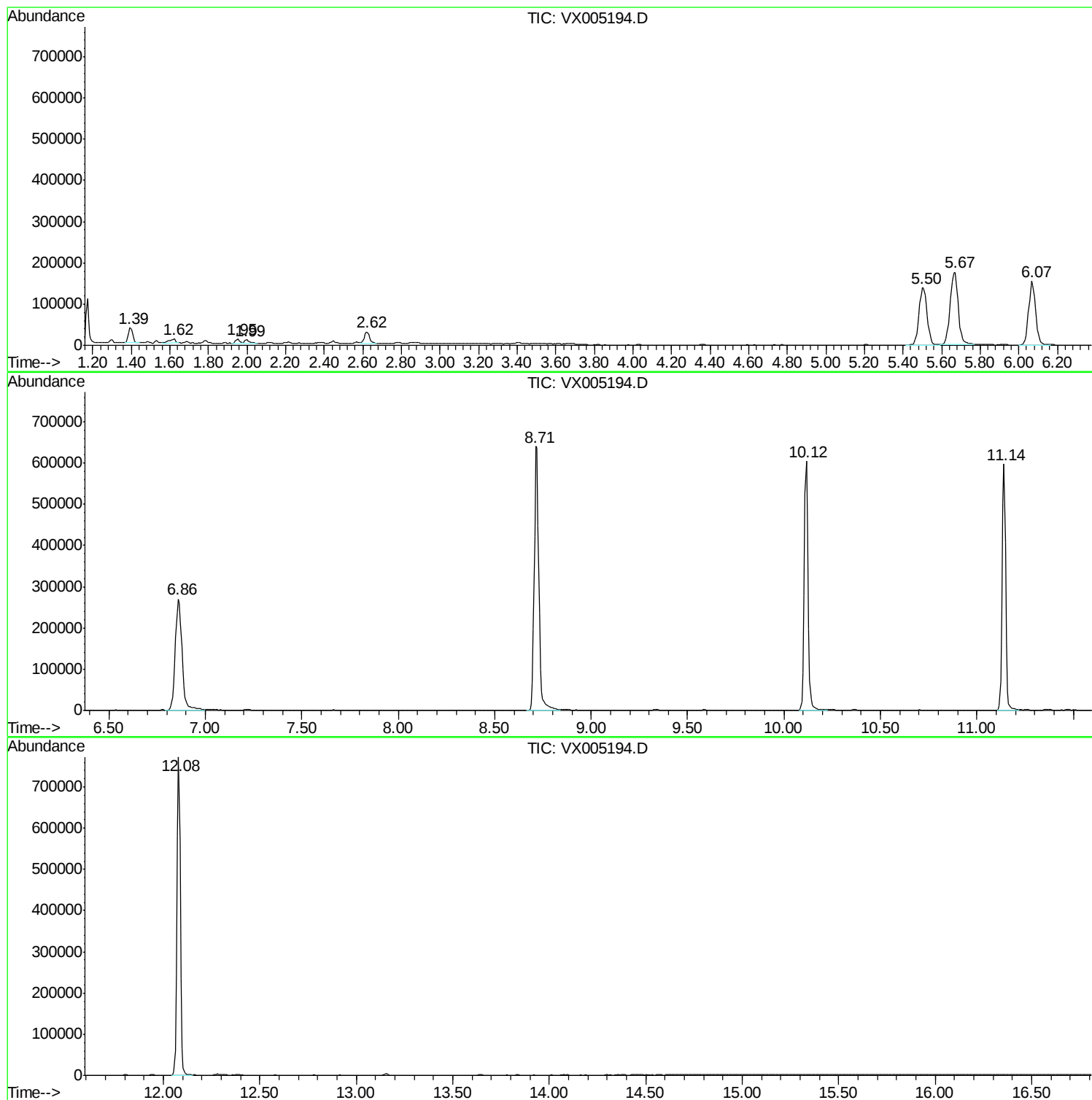
Sum of corrected areas: 5750840

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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X092618W.M
Quant Title : SW846 8260

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



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Instrument :
MSVOA_X
ClientSampled :
SA-TW514-4044

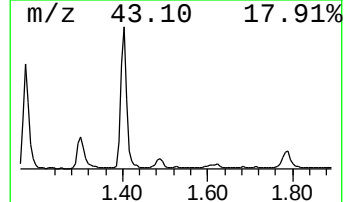
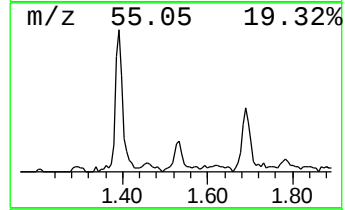
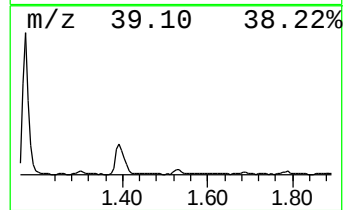
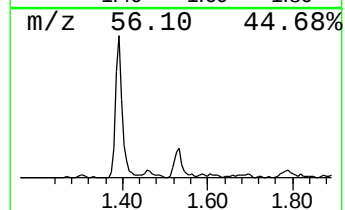
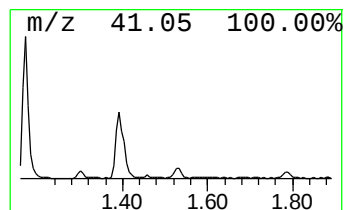
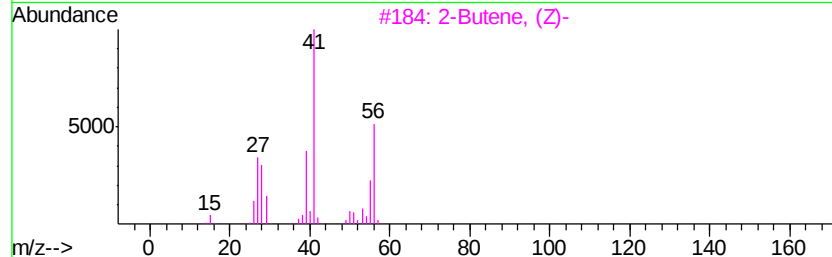
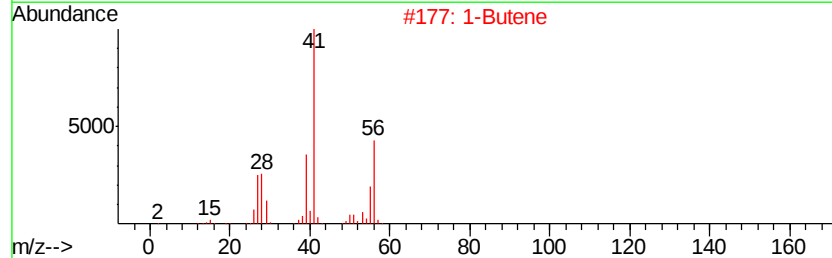
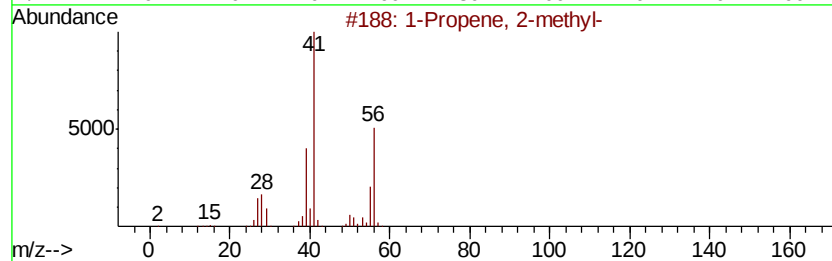
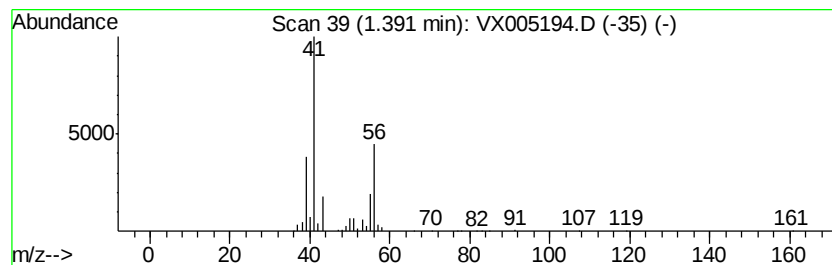
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Quant Title : SW846 8260

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.39	5.61 ug/l	57740	Pentafluorobenzene	5.67

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			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	83
4			2-Butene, (E)-	56	C4H8	000624-64-6	80
5			2-Butene	56	C4H8	000107-01-7	72



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
1-Propene, 2-methyl-	1.39	5.6	ug/l	57740	1	5.67	514212	50.0