

Data Path : W:\HPCHEM1\MSVOA X\DATA\VX051518\
 Data File : VX001745.D
 Acq On : 15 May 2018 15:08
 Operator : JC/MD
 Sample : J2737-05
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 KY062MW140-180503

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 : W:\HPCHEM1\MSVOA_X\METHOD\82X051518W.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	1.075	76	80	92	rBV	1276561	1599810	100.00%	18.938%
2	1.172	92	96	107	rVB	134147	140840	8.80%	1.667%
3	1.386	127	131	140	rBV3	60513	99915	6.25%	1.183%
4	2.629	329	335	343	rVB2	14350	30406	1.90%	0.360%
5	2.812	359	365	370	rVB	10660	20696	1.29%	0.245%
6	4.605	648	659	671	rBV	76236	212016	13.25%	2.510%
7	5.513	795	808	823	rBV	185300	524400	32.78%	6.208%
8	5.672	823	834	851	rVV	230652	653802	40.87%	7.740%
9	6.074	889	900	909	rBV	196912	518000	32.38%	6.132%
10	6.488	959	968	979	rVB2	13524	35894	2.24%	0.425%
11	6.866	1021	1030	1044	rBV	319795	759068	47.45%	8.986%
12	8.720	1327	1334	1343	rBV	840064	1331498	83.23%	15.762%
13	9.342	1426	1436	1444	rVB4	8986	16860	1.05%	0.200%
14	10.116	1555	1563	1575	rBV	571873	812851	50.81%	9.623%
15	11.146	1726	1732	1746	rBV	656415	864419	54.03%	10.233%
16	12.085	1880	1886	1898	rBV	671408	826923	51.69%	9.789%

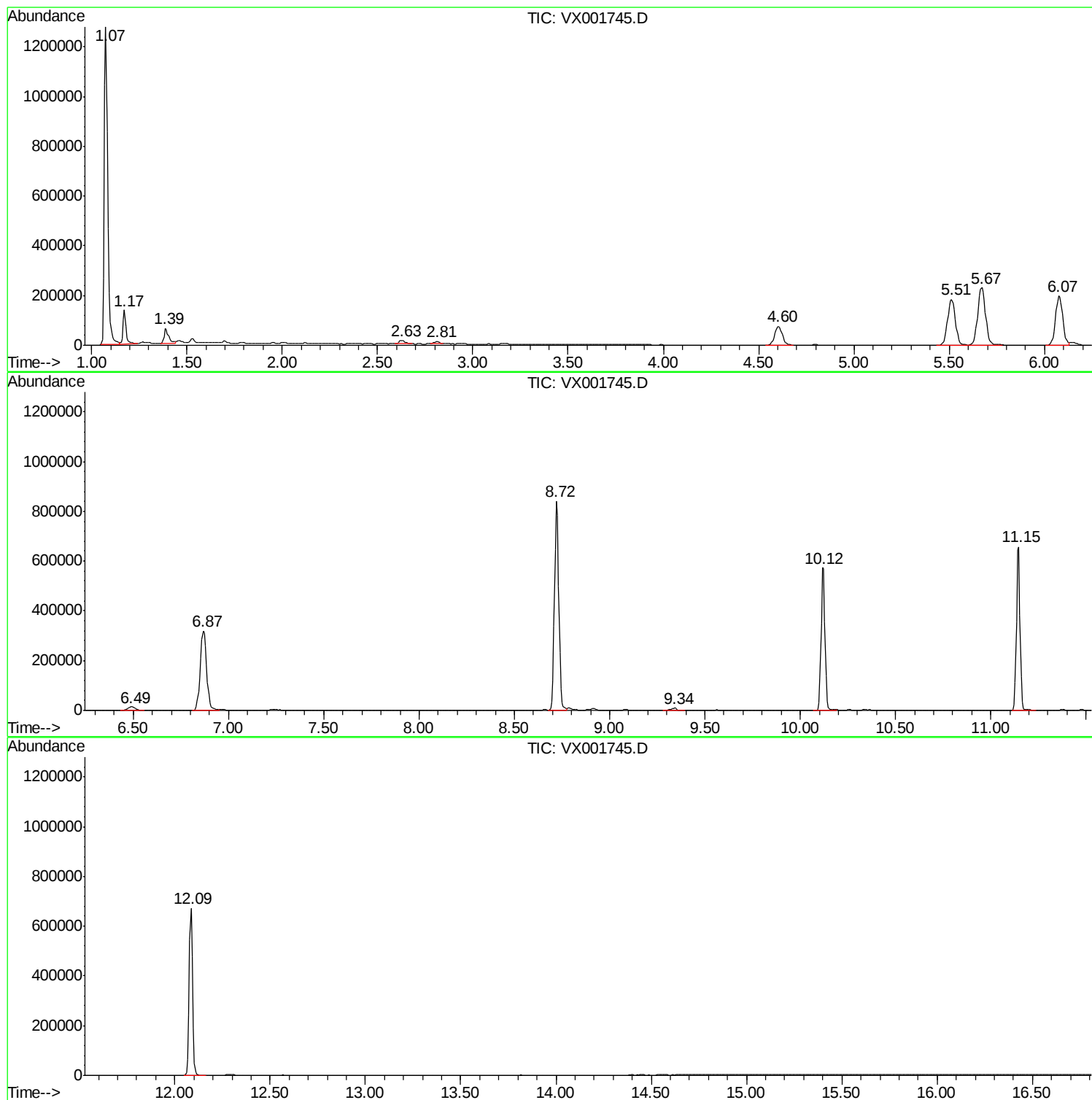
Sum of corrected areas: 8447398

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Quant Title : SW846 8260

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



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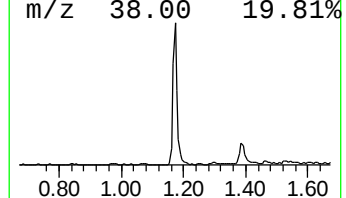
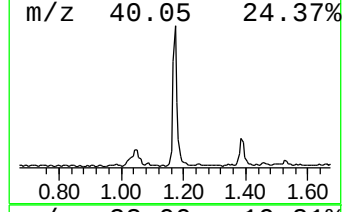
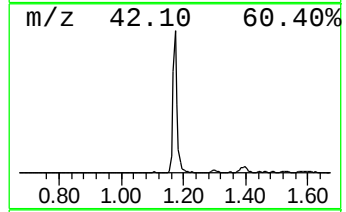
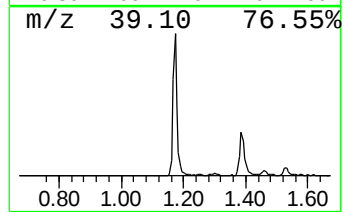
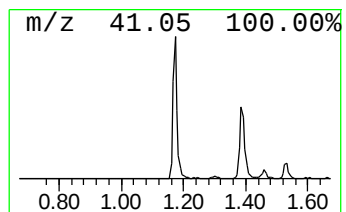
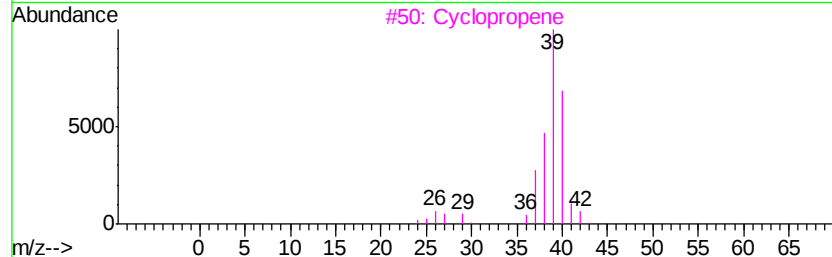
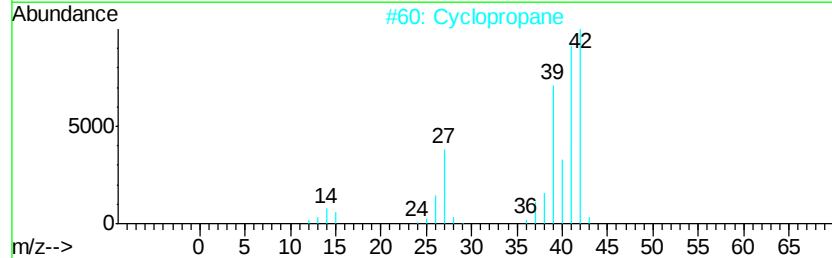
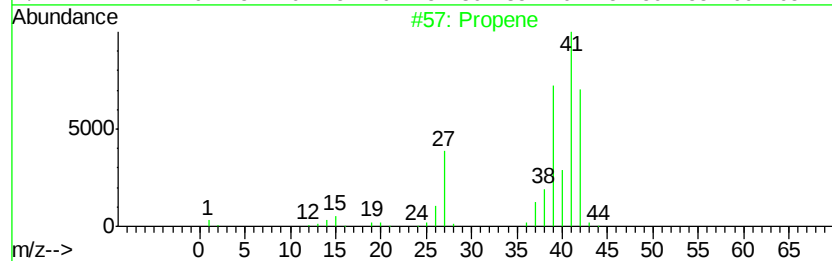
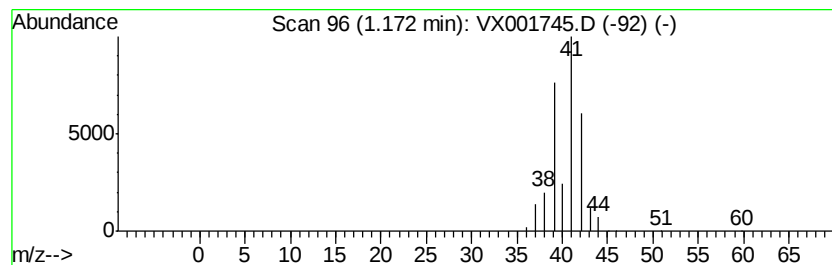
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Peak Number 2 Propene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.17	10.77 ug/l	140840	Pentafluorobenzene	5.67

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propene	42	C3H6	000115-07-1	90
2			Cyclopropane	42	C3H6	000075-19-4	9
3			Cyclopropene	40	C3H4	002781-85-3	9
4			2-Oxetanone, 4-methyl-	86	C4H6O2	003068-88-0	4
5			Acetonitrile	41	C2H3N	000075-05-8	3



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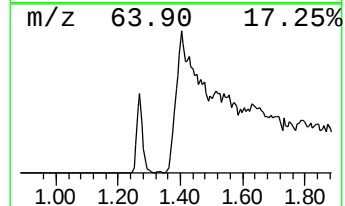
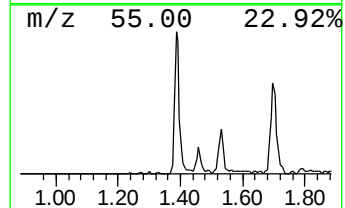
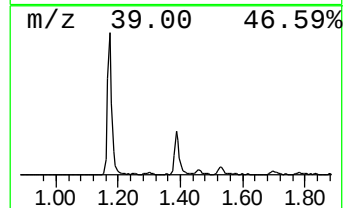
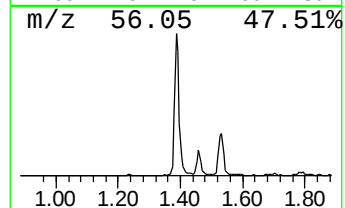
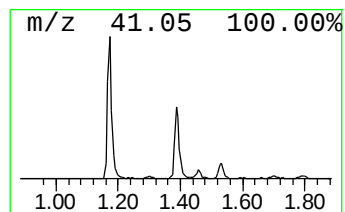
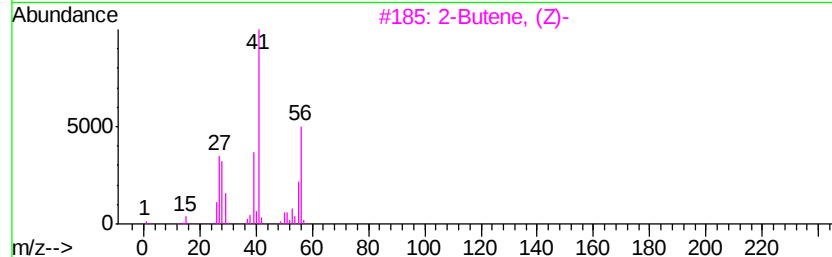
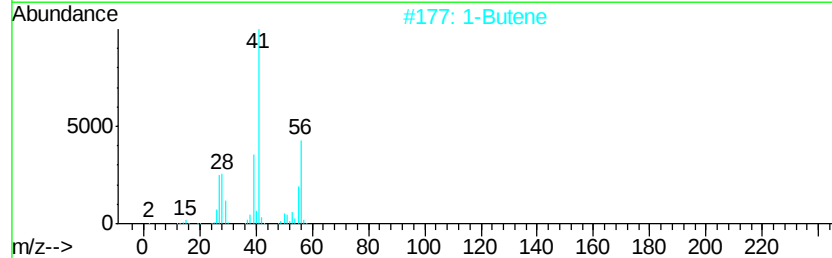
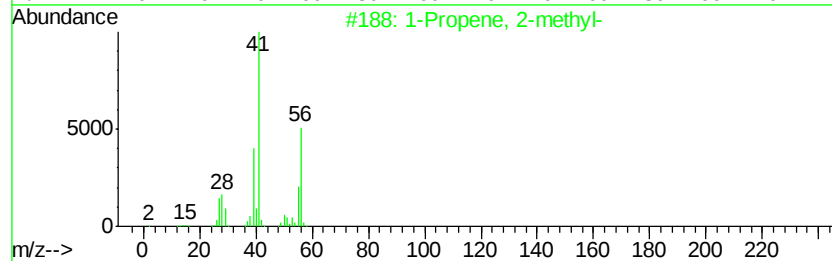
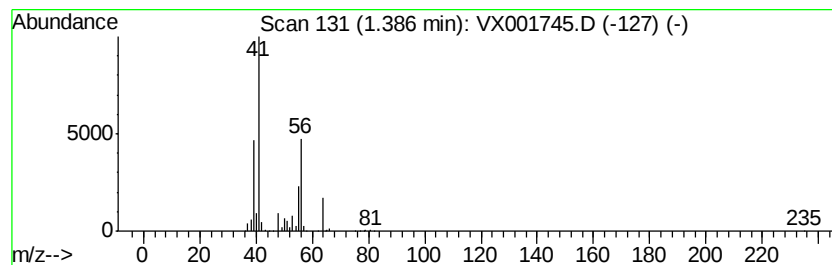
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Peak Number 3 1-Propene, 2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.39	7.64 ug/l	99915	Pentafluorobenzene	5.67

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	1.17	10.8	ug/l	140840	1	5.67	653802	50.0
1-Propene, 2-methyl-	1.39	7.6	ug/l	99915	1	5.67	653802	50.0