

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX060221\
 Data File : VX022338.D
 Acq On : 02 Jun 2021 15:04
 Operator : JC/MD
 Sample : M2572-12
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 TB03-20210527

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\82X051921W.M

Title : SW846 8260

Signal : TIC: VX022338.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.270	24	30	36	rBV3	9046	11658	2.61%	0.504%
2	1.355	41	44	51	rVB	45136	41075	9.18%	1.777%
3	1.453	57	60	67	rBV	9131	10374	2.32%	0.449%
4	1.569	70	79	82	rVB3	2968	5027	1.12%	0.218%
5	2.386	207	213	221	rBV2	5077	8301	1.86%	0.359%
6	2.550	234	240	246	rVB	4554	8457	1.89%	0.366%
7	2.965	300	308	318	rBV6	2154	5908	1.32%	0.256%
8	4.574	565	572	585	rVB3	3038	10078	2.25%	0.436%
9	5.397	694	707	719	rBV2	46725	132365	29.60%	5.727%
10	5.562	723	734	748	rBV	72074	205827	46.02%	8.906%
11	5.970	787	801	811	rBV2	59131	156954	35.09%	6.791%
12	6.769	922	932	944	rBV	132771	315293	70.50%	13.643%
13	8.653	1234	1241	1255	rBV	287245	447247	100.00%	19.352%
14	10.055	1465	1471	1482	rBV	263445	371736	83.12%	16.085%
15	11.085	1635	1640	1651	rVB	206792	263375	58.89%	11.396%
16	12.024	1789	1794	1804	rVB	257000	317430	70.97%	13.735%

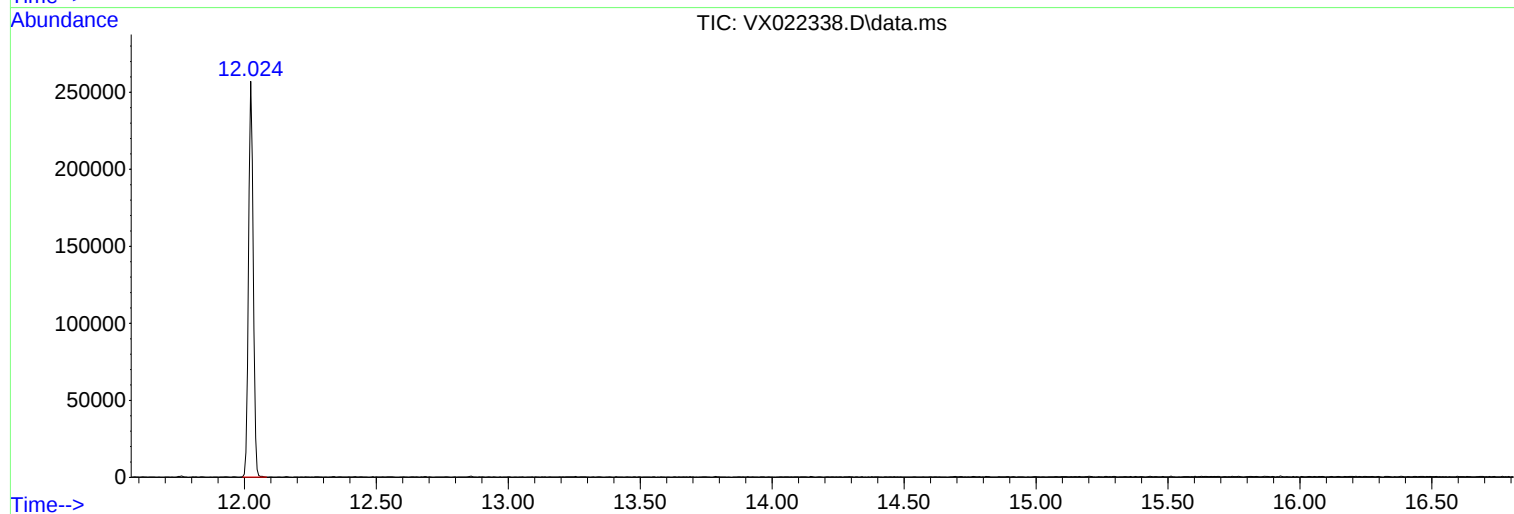
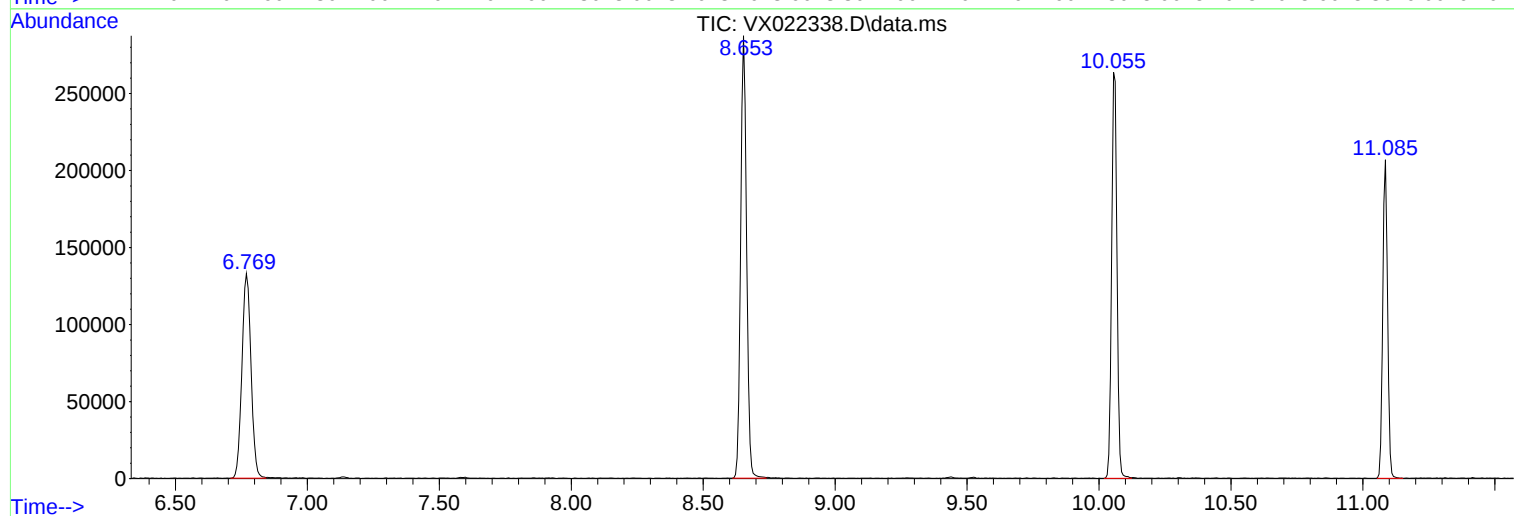
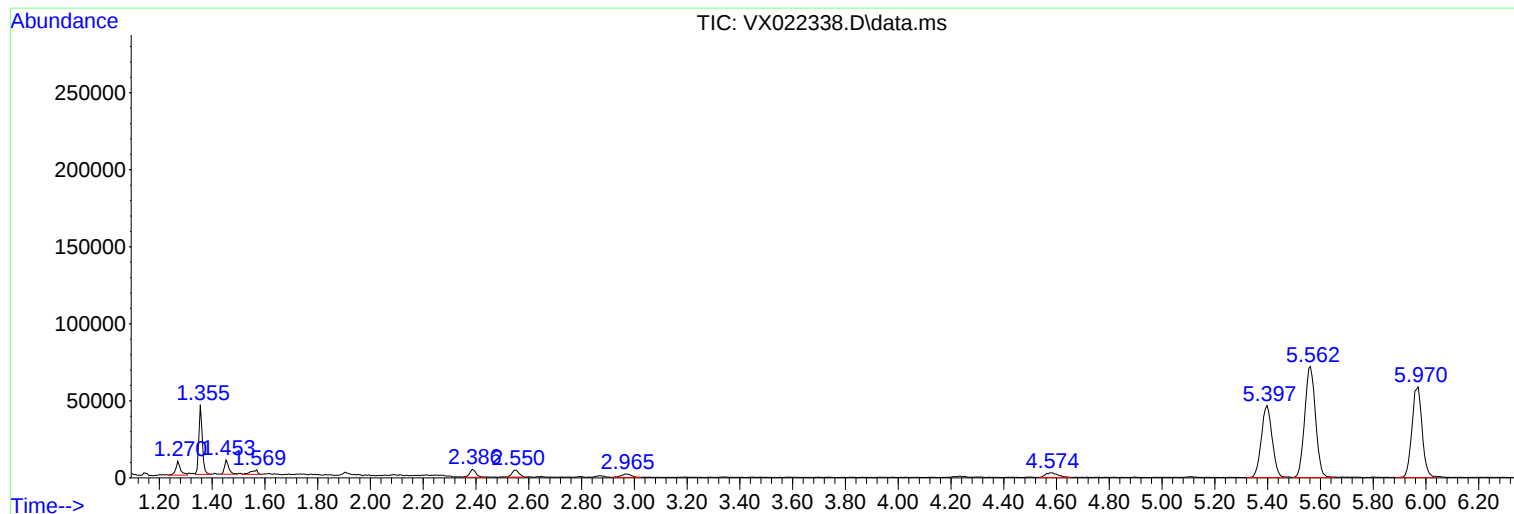
Sum of corrected areas: 2311105

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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X051921W.M
Quant Title : SW846 8260

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



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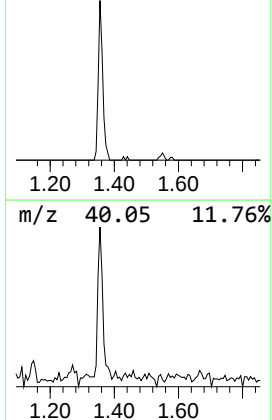
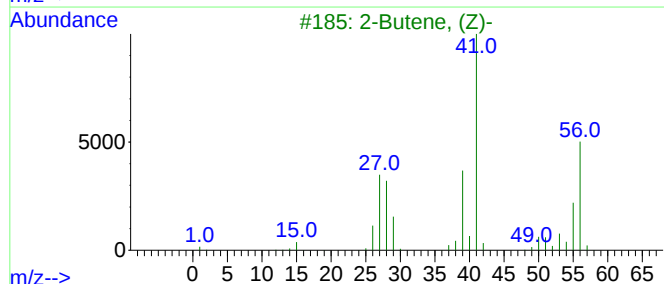
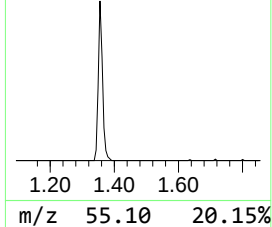
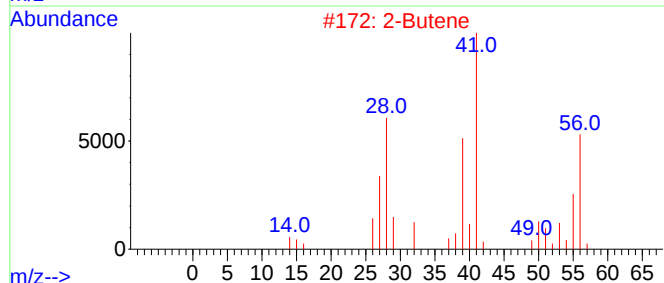
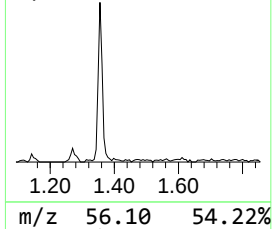
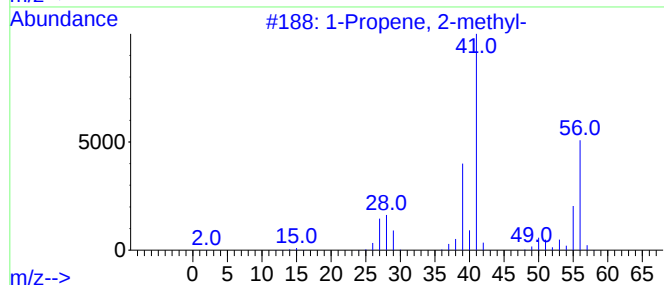
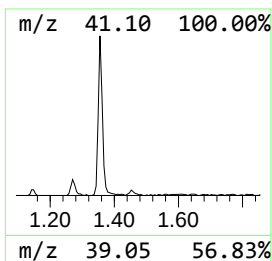
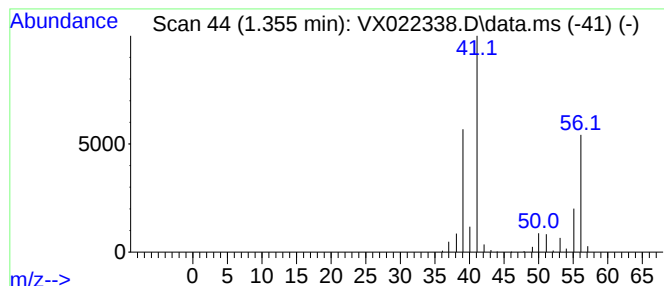
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Peak Number 1 1-Propene, 2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.355	9.98 ug/l	41075	Pentafluorobenzene	5.556

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Propene, 2-methyl-			56	C4H8	000115-11-7	87
2	2-Butene			56	C4H8	000107-01-7	80
3	2-Butene, (Z)-			56	C4H8	000590-18-1	50
4	2-Butene, (E)-			56	C4H8	000624-64-6	50
5	1-Butene			56	C4H8	000106-98-9	49



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
1-Propene, 2-me...	1.355	10.0	ug/l	41075	1	5.556	205827	50.0