

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX072722\
 Data File : VX030228.D
 Acq On : 27 Jul 2022 13:54
 Operator : JC/MD
 Sample : N3876-03
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 SPC-11-84

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\82X072522W.M

Title : SW846 8260

Signal : TIC: VX030228.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.142	6	9	16	rVB	78458	66220	5.95%	1.271%
2	1.356	41	44	53	rVB2	21100	33950	3.05%	0.651%
3	5.391	694	706	720	rBV2	124321	369225	33.20%	7.085%
4	5.556	722	733	748	rVB	145997	405130	36.43%	7.774%
5	5.964	789	800	812	rBV	150781	407988	36.68%	7.829%
6	6.769	923	932	944	rBV	233690	553016	49.72%	10.612%
7	7.135	984	992	998	rBV3	6947	14822	1.33%	0.284%
8	8.653	1234	1241	1250	rBV	720010	1112210	100.00%	21.342%
9	10.055	1464	1471	1480	rBV	547780	731844	65.80%	14.043%
10	11.085	1634	1640	1648	rBV	525654	678397	61.00%	13.018%
11	11.305	1672	1676	1682	rVB	8790	11808	1.06%	0.227%
12	12.024	1788	1794	1802	rBV	679946	826809	74.34%	15.865%

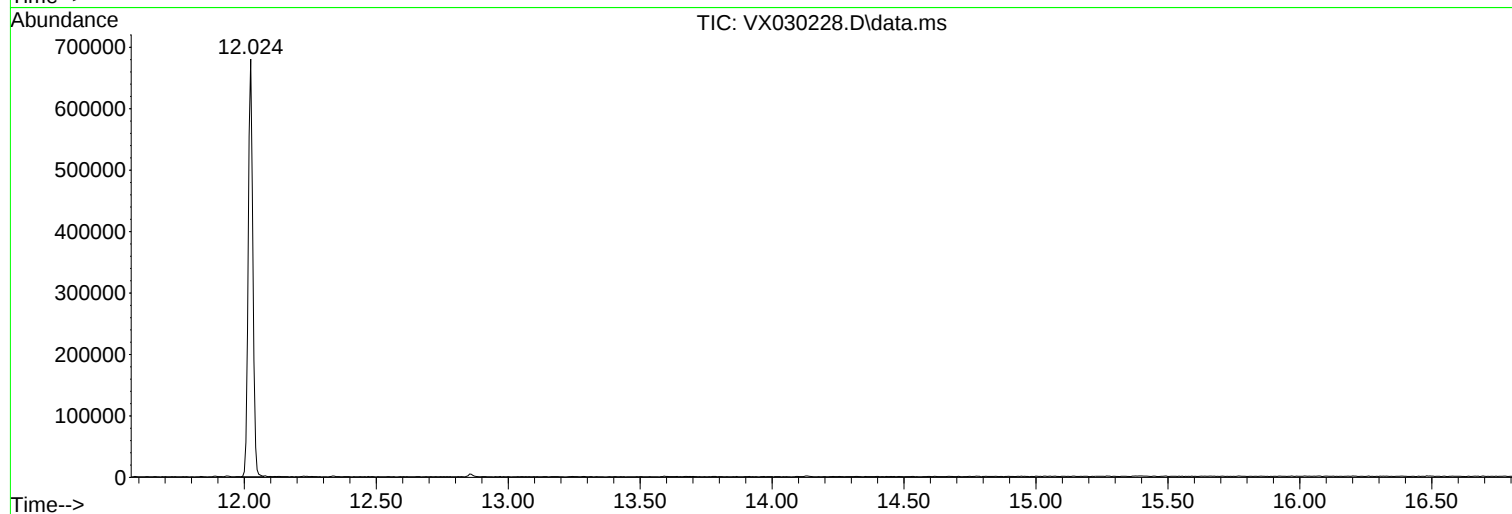
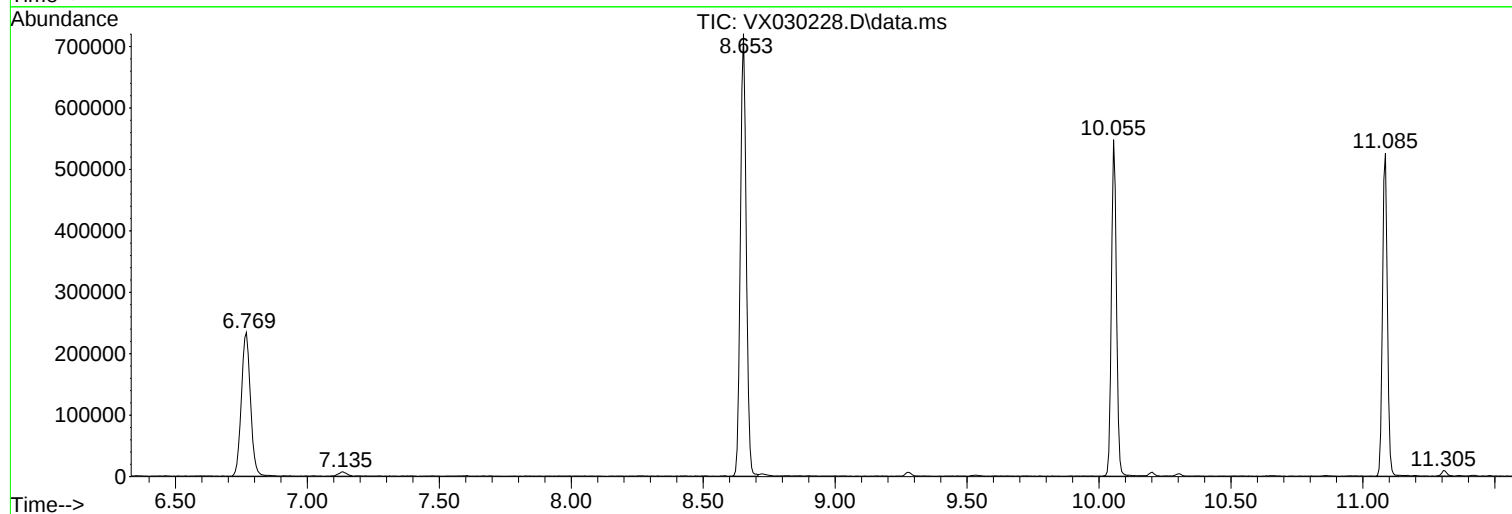
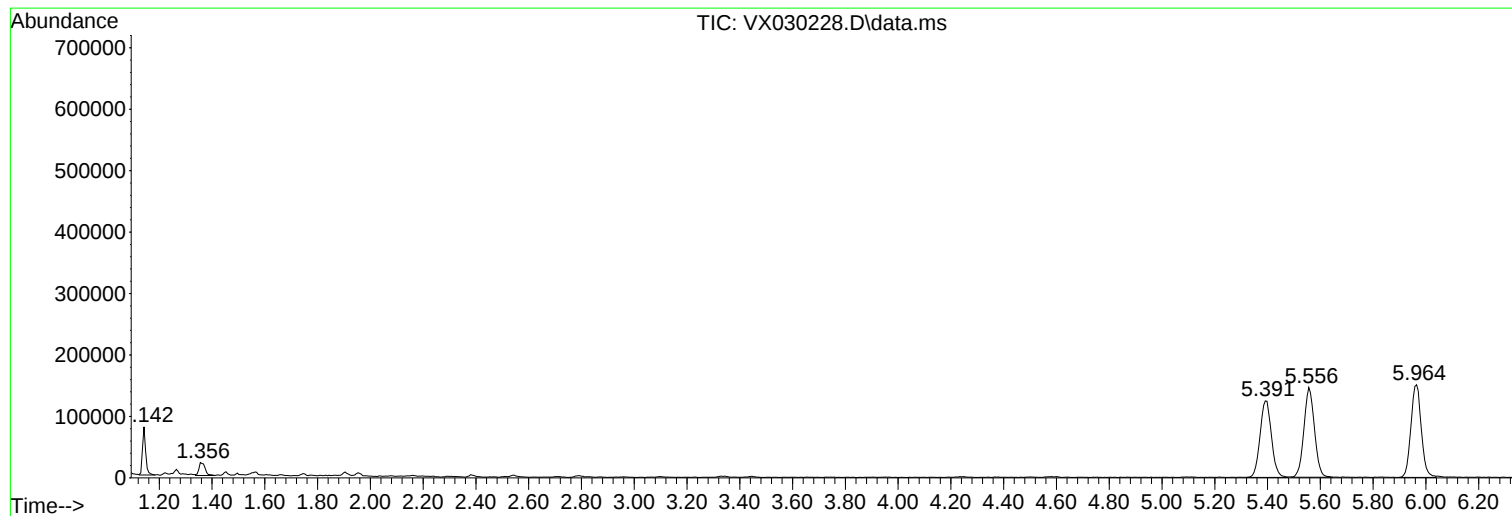
Sum of corrected areas: 5211419

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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X072522W.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P



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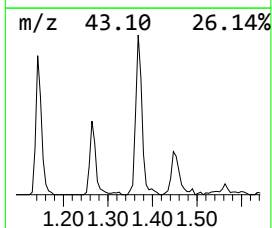
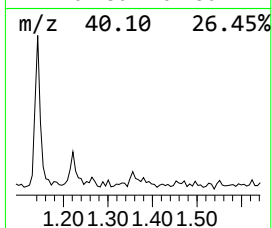
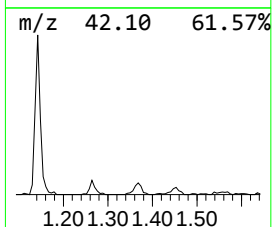
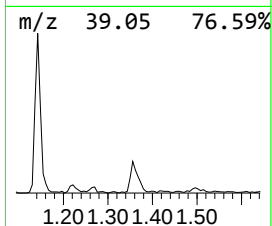
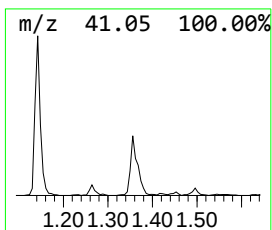
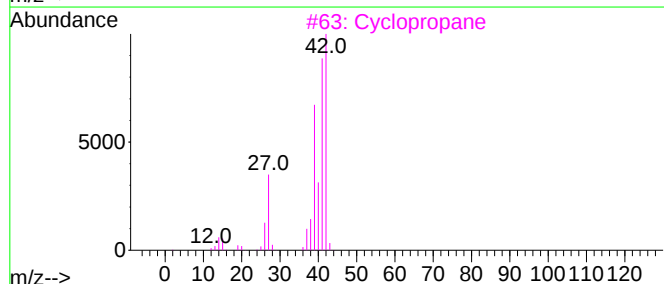
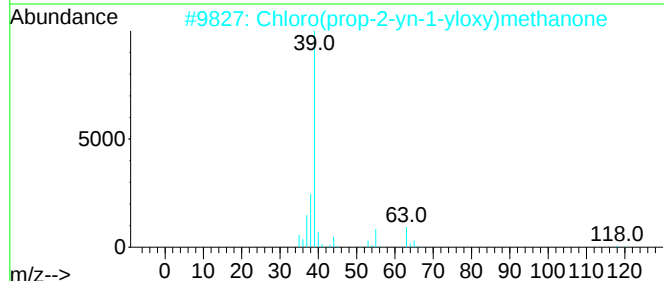
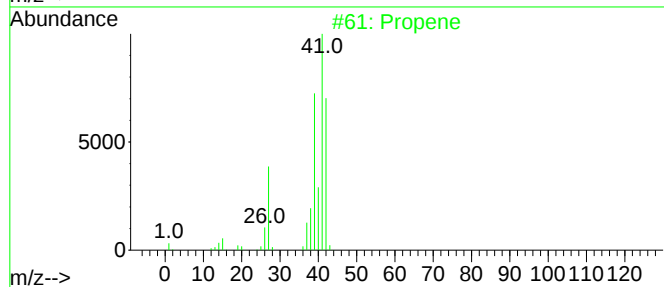
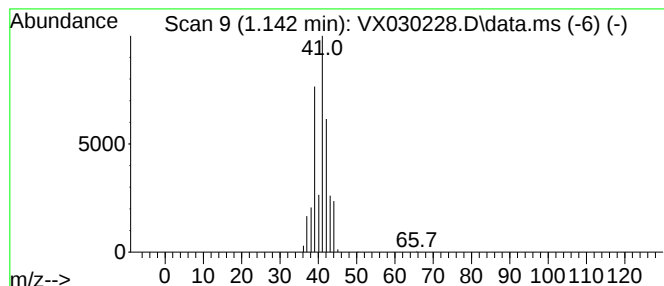
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X072522W.M
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TIC Library : C:\Database\NIST20.L
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Peak Number 1 Propene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.142	8.17 ug/l	66220	Pentafluorobenzene	5.556

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Propene	42	C3H6	000115-07-1	90
2		Chloro(prop-2-yn-1-yloxy)methanone	118	C4H3ClO2	035718-08-2	40
3		Cyclopropane	42	C3H6	000075-19-4	40
4		2-Oxetanone, 4-methyl-	86	C4H6O2	003068-88-0	9
5		Cyclopropene	40	C3H4	002781-85-3	5



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Propene	1.142	8.2	ug/l	66220	1	5.556	405130	50.0