

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY061021\
 Data File : VY005103.D
 Acq On : 11 Jun 2021 03:16
 Operator : SY/MD
 Sample : M2672-18
 Misc : 6.69G/5ML/MSVOA_Y/SOIL
 ALS Vial : 45 Sample Multiplier: 1

Instrument :
 MSVOA_Y
 ClientSampleId :
 17-B6(76-80)

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_Y\methods\82Y060221S.M

Title : SW846 8260

Signal : TIC: VY005103.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.869	4	8	26	rVB	64540	137444	10.02%	2.190%
2	2.229	61	67	76	rVB3	18002	44971	3.28%	0.717%
3	3.265	229	237	252	rVB5	8462	26635	1.94%	0.424%
4	4.729	467	477	494	rBV6	8430	32381	2.36%	0.516%
5	7.728	958	969	975	rBV2	132935	323591	23.60%	5.156%
6	7.801	975	981	992	rVB	214868	490940	35.80%	7.822%
7	8.149	1028	1038	1049	rBV	141864	312887	22.82%	4.985%
8	8.697	1119	1128	1140	rBV	365413	702213	51.21%	11.188%
9	10.185	1364	1372	1380	rBV	579772	1005560	73.33%	16.022%
10	10.941	1490	1496	1503	rBV	11224	18589	1.36%	0.296%
11	11.496	1580	1587	1597	rVB	565116	898674	65.54%	14.319%
12	12.489	1741	1750	1759	rBV2	782592	1371278	100.00%	21.849%
13	13.428	1888	1904	1910	rBV	572097	866640	63.20%	13.808%
14	13.965	1986	1992	1998	rBV	14970	25157	1.83%	0.401%
15	14.209	2026	2032	2038	rVB2	12743	19342	1.41%	0.308%

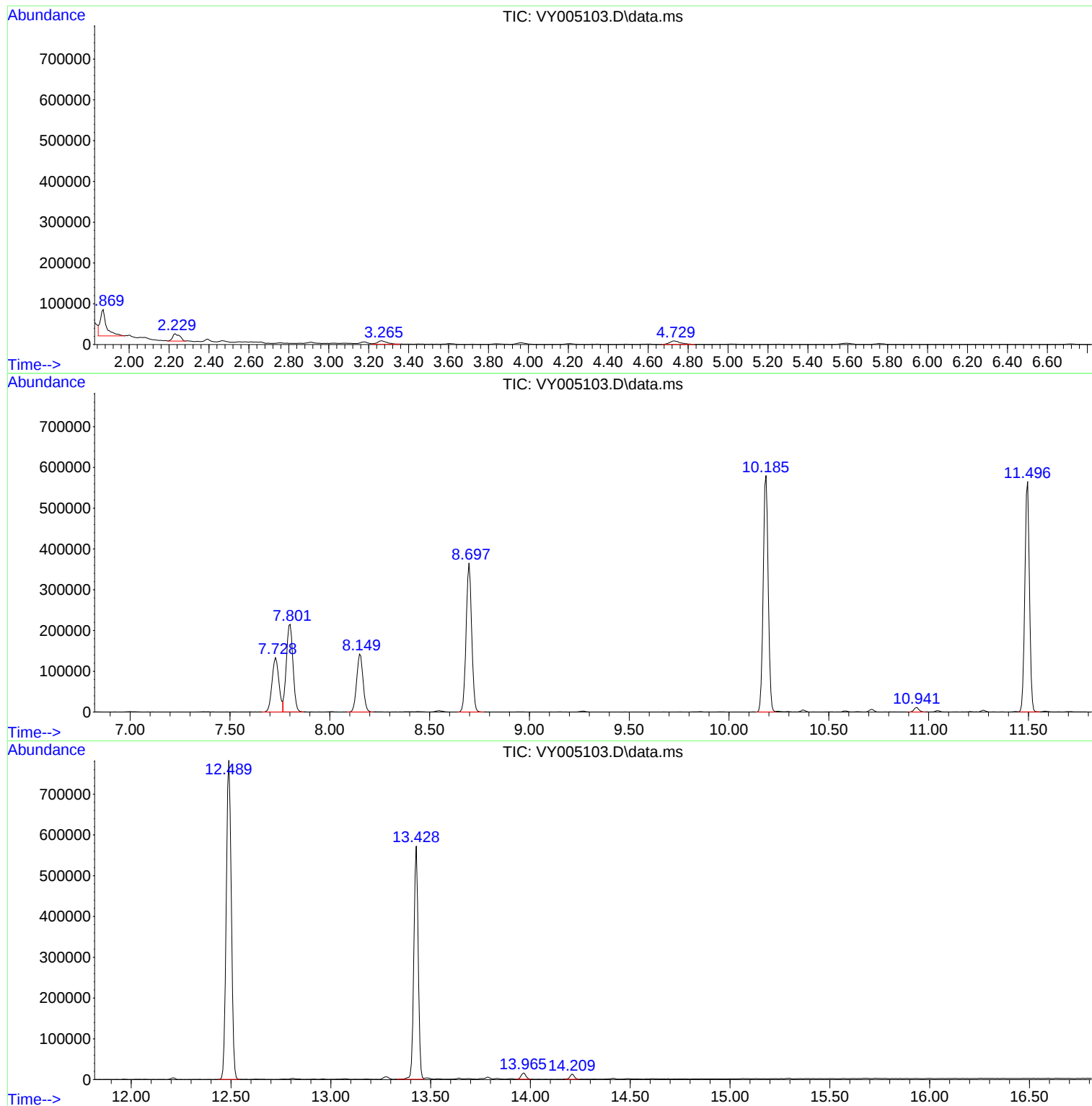
Sum of corrected areas: 6276302

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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y060221S.M
Quant Title : SW846 8260

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



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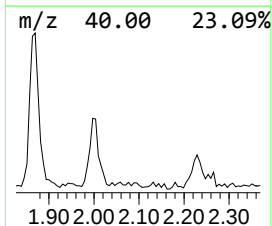
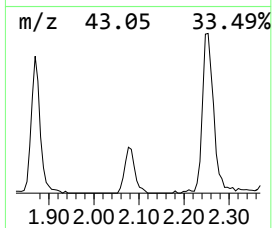
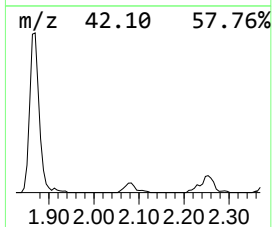
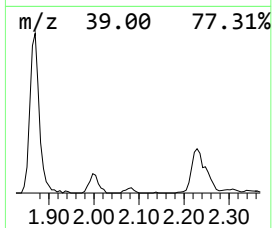
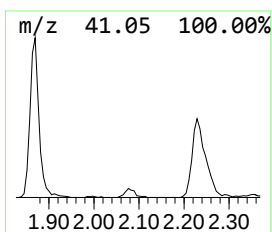
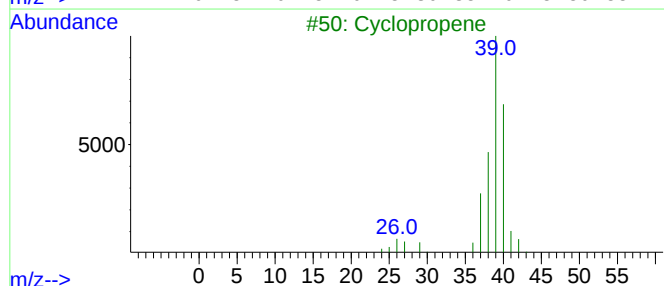
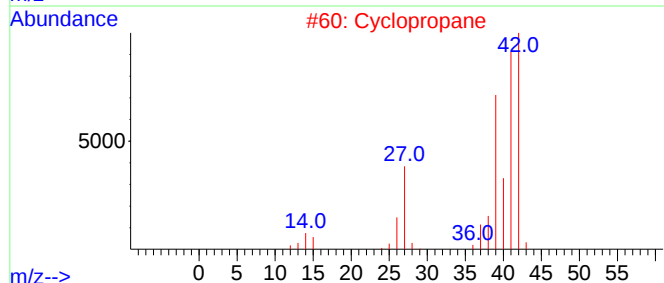
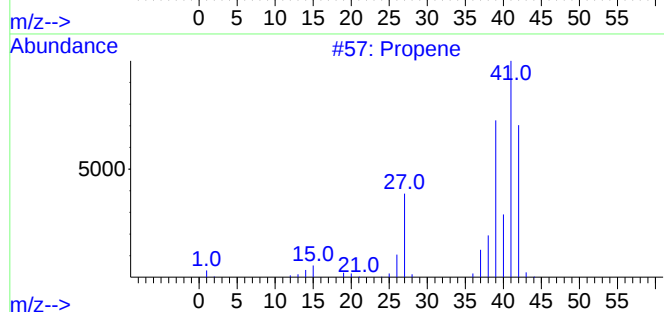
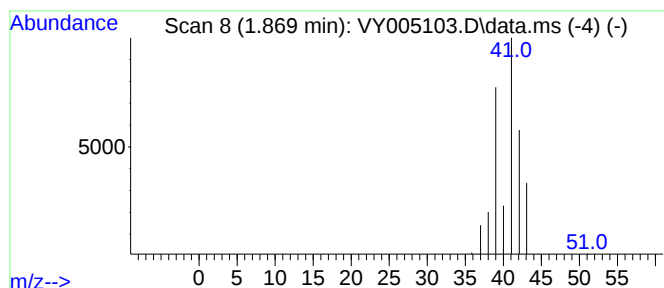
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Peak Number 1 Propene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.869	14.00 ug/l	137444	Pentafluorobenzene	7.801

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	5
4			2-Oxetanone, 4-methyl-	86	C4H6O2	003068-88-0	4
5			Acetonitrile	41	C2H3N	000075-05-8	3



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Propene	1.869	14.0	ug/l	137444	1	7.801	490940	50.0