

Data Path : Z:\VOASRV\HPCHEM1\MSVOA Y\DATA\VY050420\
 Data File : VY002569.D
 Acq On : 05 May 2020 03:05
 Operator : SY/MD
 Sample : L2475-02
 Misc : 4.07G/5ML/MSVOA Y/SOIL
 ALS Vial : 46 Sample Multiplier: 1

Instrument :
 MSVOA_Y
 ClientSampleId :
 FK-GF-03-009-A

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\82Y043020S.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	1.883	19	24	36	rVB	181153	259898	14.56%	2.684%
2	2.268	80	87	93	rVB	11063	19854	1.11%	0.205%
3	4.755	485	495	504	rBV2	18136	58424	3.27%	0.603%
4	4.846	504	510	524	rVB2	6282	19714	1.10%	0.204%
5	7.748	974	986	992	rBV	252248	602346	33.74%	6.221%
6	7.821	992	998	1014	rVB	442025	989192	55.41%	10.216%
7	8.169	1045	1055	1067	rBV	225590	498583	27.93%	5.149%
8	8.711	1136	1144	1155	rVB	644960	1293214	72.44%	13.355%
9	10.199	1380	1388	1396	rBV	1046377	1785225	100.00%	18.436%
10	11.504	1593	1602	1613	rBV	938430	1493387	83.65%	15.422%
11	12.497	1757	1765	1773	rBV	712990	1133063	63.47%	11.701%
12	13.436	1912	1919	1927	rBV	908445	1402206	78.55%	14.481%
13	13.540	1930	1936	1946	rBV	55705	89646	5.02%	0.926%
14	13.711	1957	1964	1967	rBV6	8275	19350	1.08%	0.200%
15	13.979	2003	2008	2014	rVB2	11360	19091	1.07%	0.197%

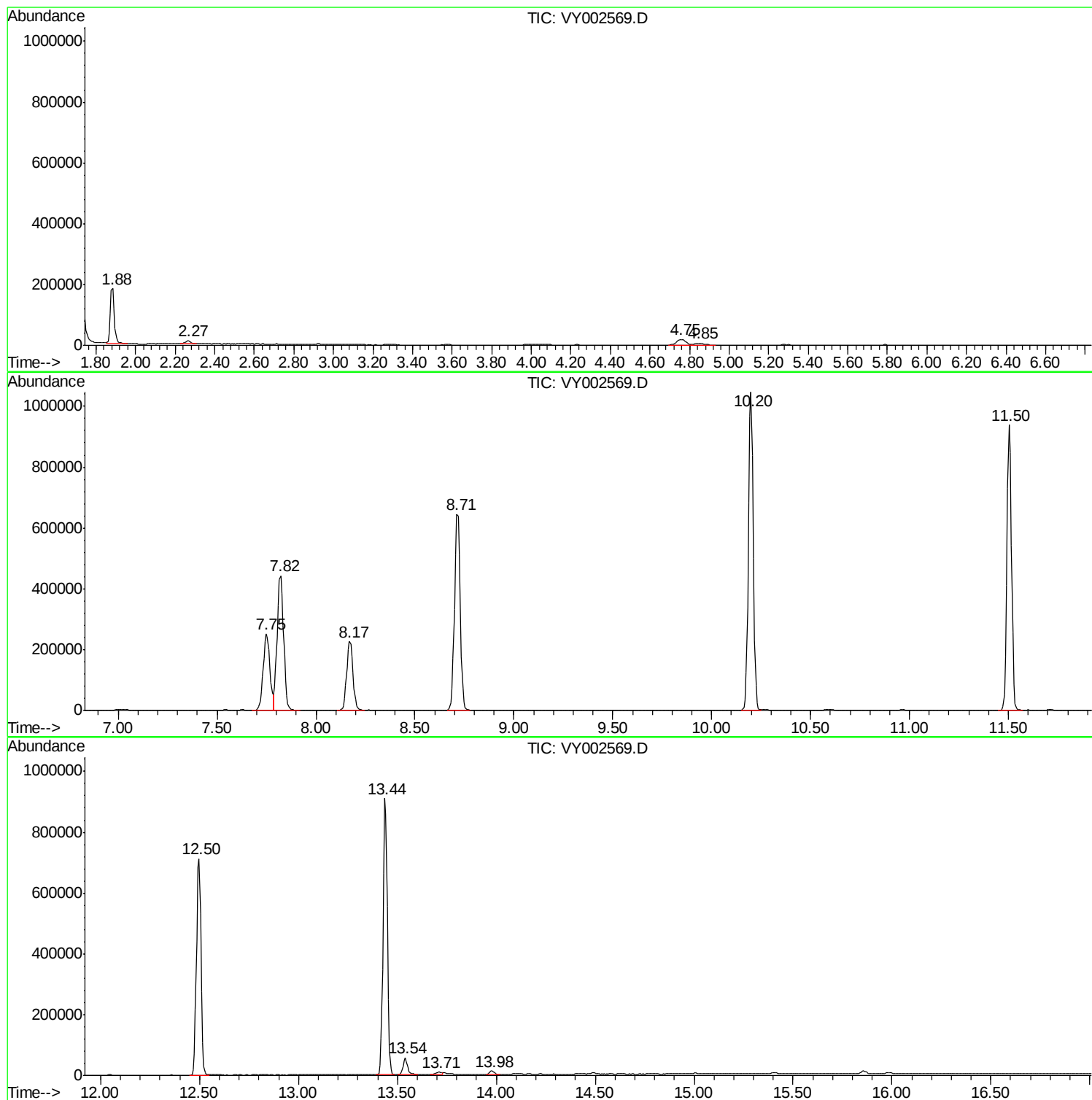
Sum of corrected areas: 9683193

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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_Y\METHODS\82Y043020S.M
Quant Title : SW846 8260

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



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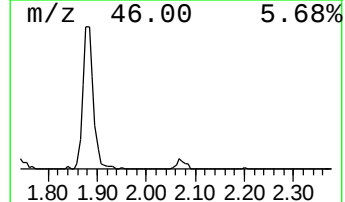
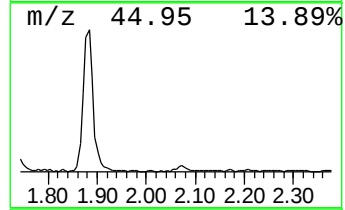
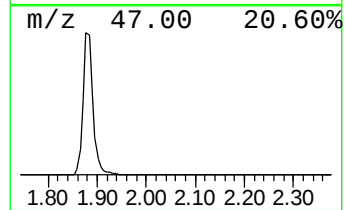
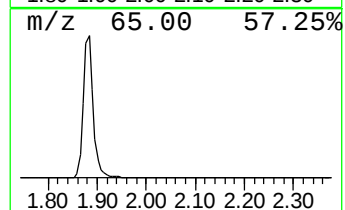
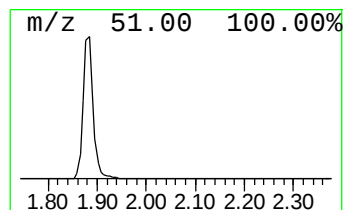
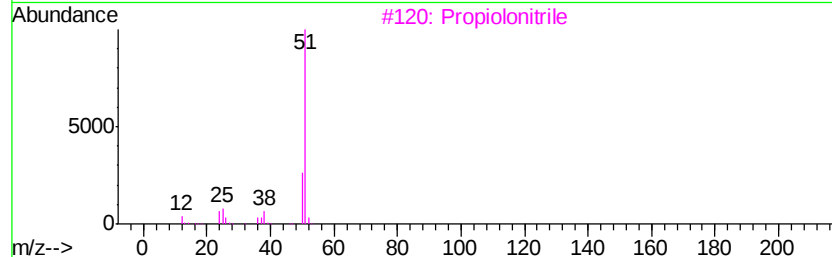
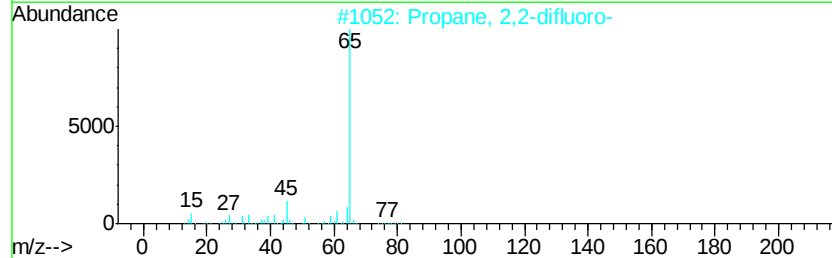
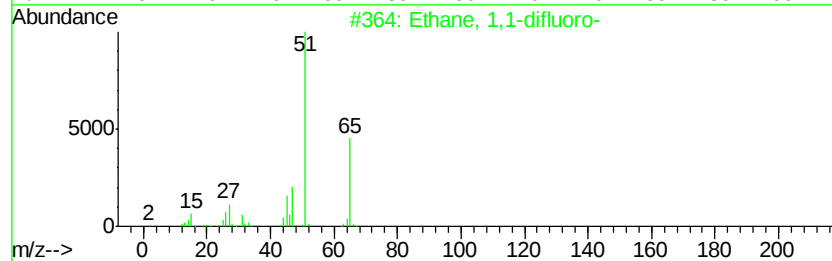
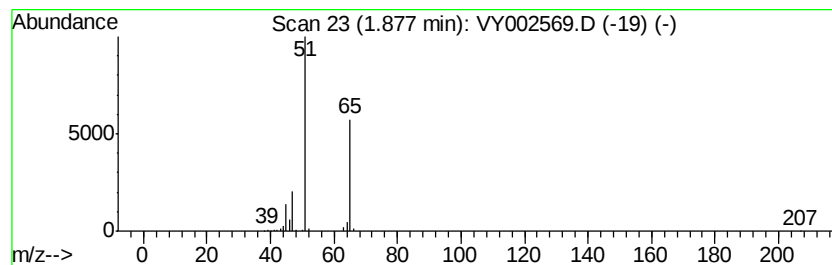
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Peak Number 1 Ethane, 1,1-difluoro- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.88	13.14 ug/l	259898	Pentafluorobenzene	7.82

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Ethane, 1,1-difluoro-	66	C2H4F2	000075-37-6	91
2		Propane, 2,2-difluoro-	80	C3H6F2	000420-45-1	9
3		Propiolonitrile	51	C3HN	001070-71-9	3
4		Difluoromethane	52	CH2F2	000075-10-5	1
5		2-Chloroethanol	80	C2H5ClO	000107-07-3	1



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Ethane, 1,1-diflu...	1.88	13.1	ug/l	259898	1	7.82	989192	50.0