

Data Path : Z:\VOASRV\HPCHEM1\MSVOA Y\DATA\VY013120\
 Data File : VY001460.D
 Acq On : 31 Jan 2020 17:00
 Operator : SY/MD
 Sample : L1333-01
 Misc : 7.00G/5ML/MSVOA Y/SOIL
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_Y
 ClientSampleId :
 TP-1-2

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 : Z:\VOASRV\HPCHEM1\MSVOA_Y\METHODS\82Y012120S.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	1.871	17	22	37	rVB	87773	122767	13.32%	2.406%
2	2.255	77	85	92	rVB2	6192	11541	1.25%	0.226%
3	4.749	481	494	504	rBV3	9900	33141	3.60%	0.650%
4	7.742	976	985	991	rBV	118157	283226	30.73%	5.552%
5	7.815	991	997	1007	rVB	210304	460948	50.01%	9.035%
6	8.163	1044	1054	1067	rVB	124037	274890	29.82%	5.388%
7	8.712	1134	1144	1159	rBV	336893	671790	72.89%	13.168%
8	10.193	1380	1387	1395	rBV	534226	921697	100.00%	18.066%
9	10.589	1445	1452	1461	rBV3	7616	15158	1.64%	0.297%
10	11.504	1595	1602	1609	rBV	497382	812122	88.11%	15.919%
11	12.491	1757	1764	1772	rBV	385091	633648	68.75%	12.420%
12	13.132	1863	1869	1874	rBV9	5066	11132	1.21%	0.218%
13	13.436	1913	1919	1926	rBV	500554	759679	82.42%	14.891%
14	13.540	1933	1936	1944	rVB5	6092	10884	1.18%	0.213%
15	13.973	2001	2007	2013	rVB2	19388	33867	3.67%	0.664%
16	14.418	2073	2080	2086	rBV3	7107	15769	1.71%	0.309%
17	15.399	2237	2241	2246	rVB	18235	29468	3.20%	0.578%

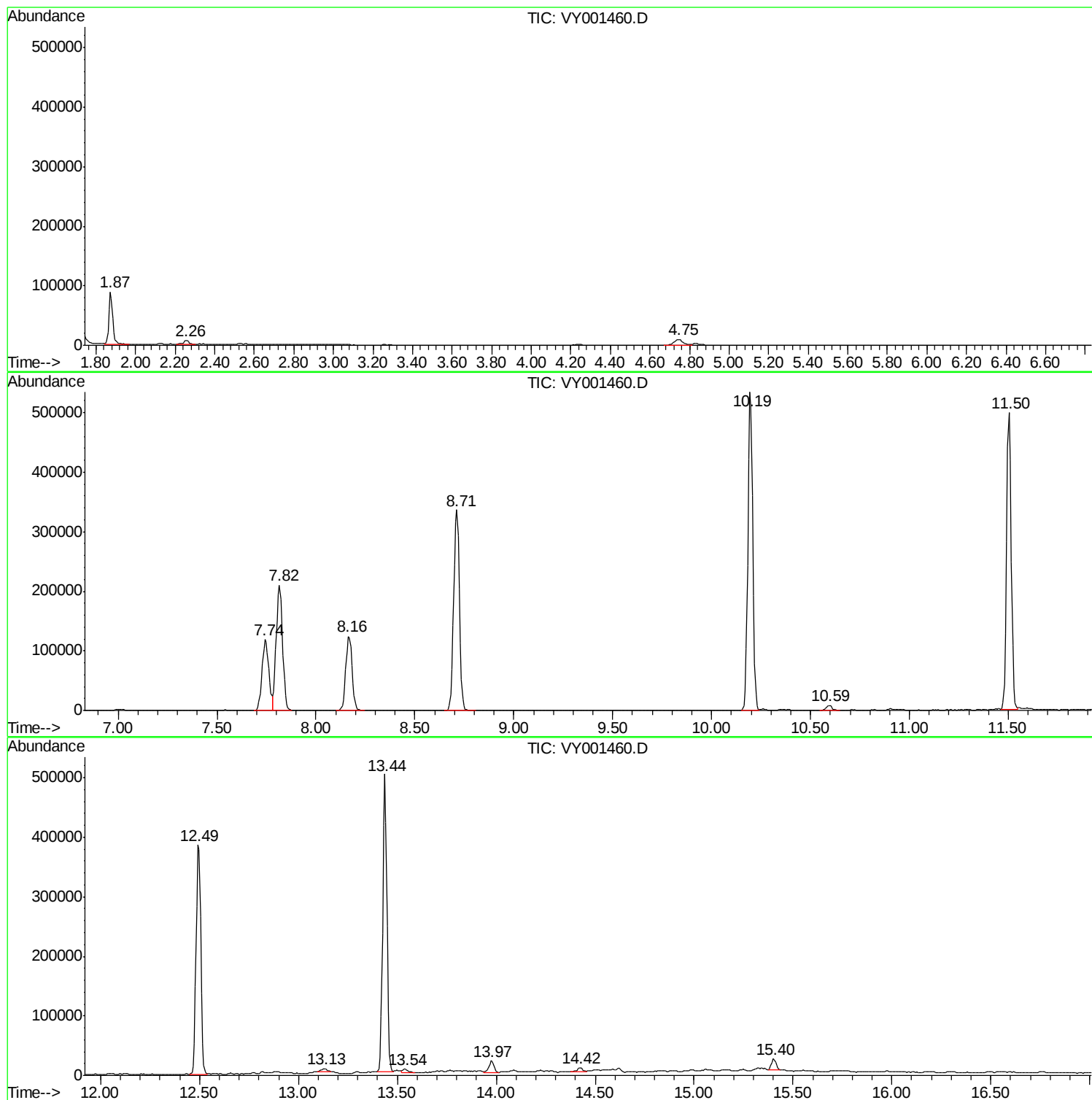
Sum of corrected areas: 5101727

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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_Y\METHODS\82Y012120S.M
Quant Title : SW846 8260

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



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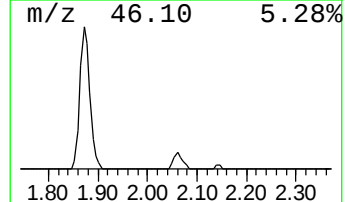
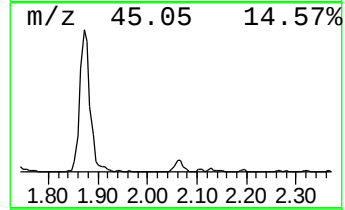
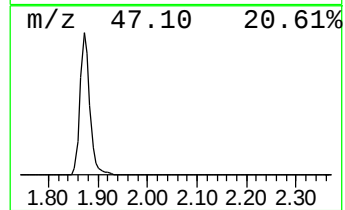
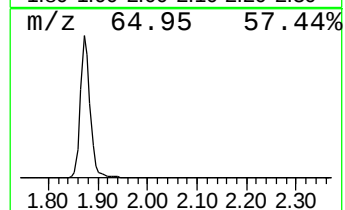
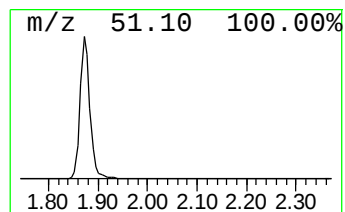
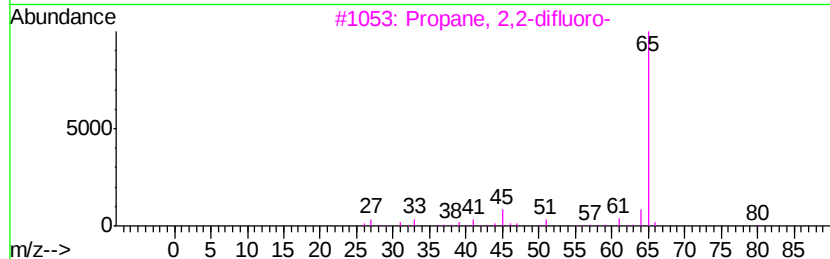
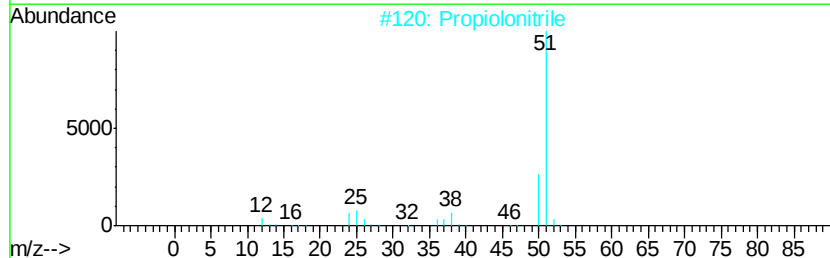
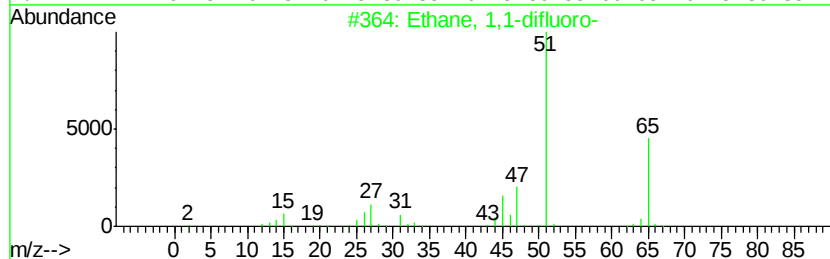
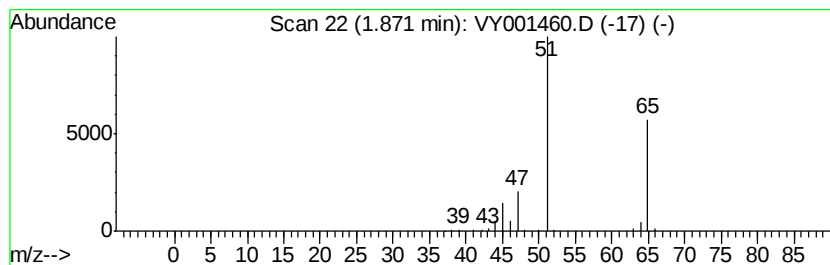
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Peak Number 1 Ethane, 1,1-difluoro- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.87	13.32 ug/l	122767	Pentafluorobenzene	7.82

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Ethane, 1,1-difluoro-	66	C2H4F2	000075-37-6	91
2		Propiolonitrile	51	C3HN	001070-71-9	3
3		Propane, 2,2-difluoro-	80	C3H6F2	000420-45-1	2
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...	1.87	13.3	ug/l	122767	1	7.82	460948	50.0