

Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD012221\
 Data File : VD068155.D
 Acq On : 22 Jan 2021 15:04
 Operator : VA/SY
 Sample : M1209-02
 Misc : 5.5G/5.00ml/MSVOA_D/SOIL
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 FK-PS-01-069

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_D\Method\82D011921S.M
 Title : SW846 8260

Signal : TIC: VD068155.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.950	3	5	14	rVB	426257	558836	18.19%	3.577%
2	7.920	1009	1020	1025	rBV2	377857	918458	29.90%	5.879%
3	7.985	1025	1031	1043	rVB	733962	1635098	53.22%	10.466%
4	8.338	1080	1091	1103	rBV	337126	790637	25.74%	5.061%
5	8.867	1172	1181	1196	rBV	1040769	2166561	70.52%	13.868%
6	10.344	1421	1432	1449	rBV	1677883	3072102	100.00%	19.665%
7	11.649	1646	1654	1666	rVB	1379459	2448826	79.71%	15.675%
8	12.632	1811	1821	1831	rBV	1056284	1772137	57.68%	11.344%
9	13.573	1974	1981	1991	rVB	1266262	2127488	69.25%	13.618%
10	14.096	2065	2070	2077	rVB	32344	51985	1.69%	0.333%
11	15.543	2309	2316	2323	rBV2	40400	80296	2.61%	0.514%

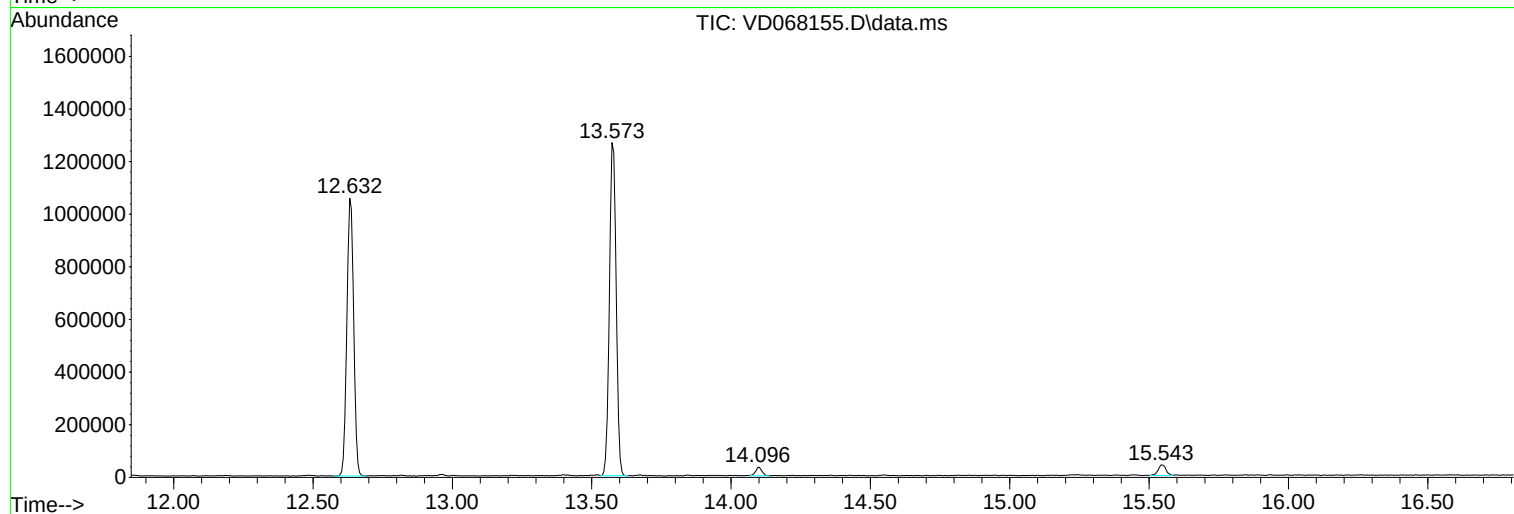
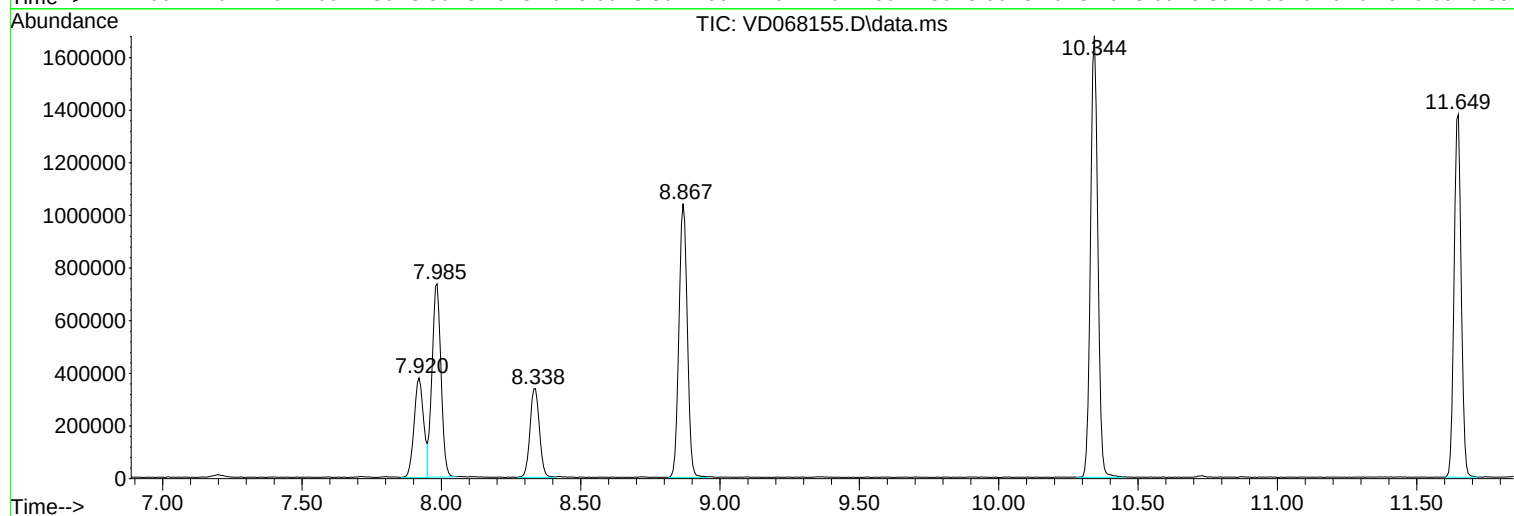
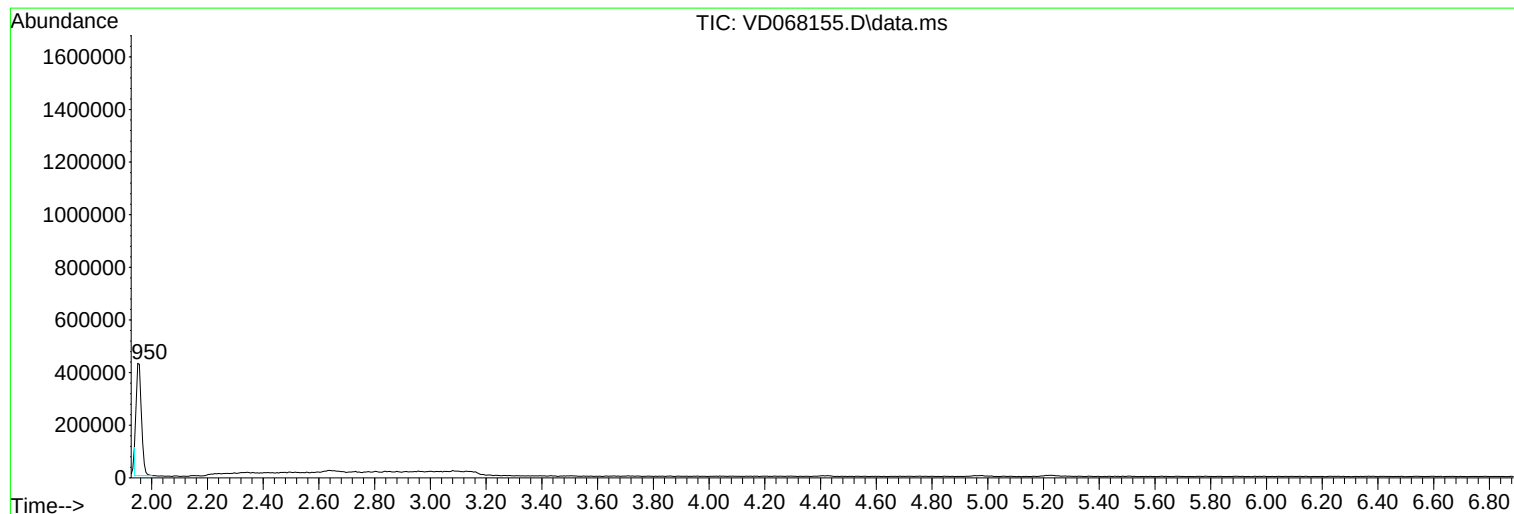
Sum of corrected areas: 15622424

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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D011921S.M
Quant Title : SW846 8260

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



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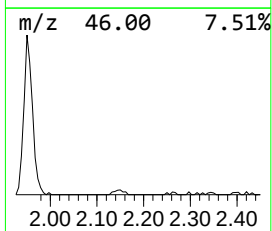
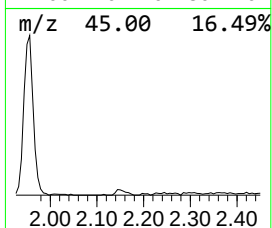
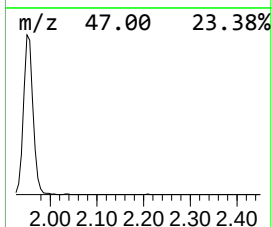
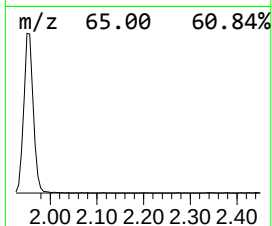
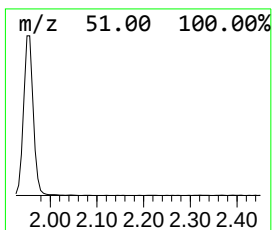
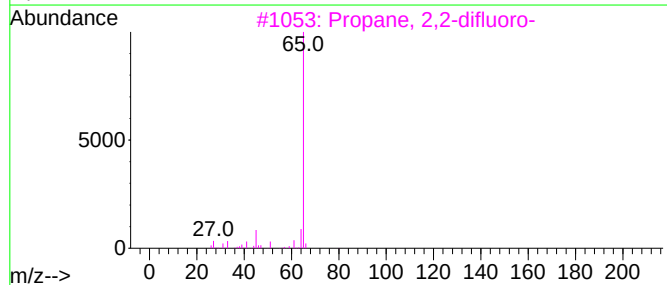
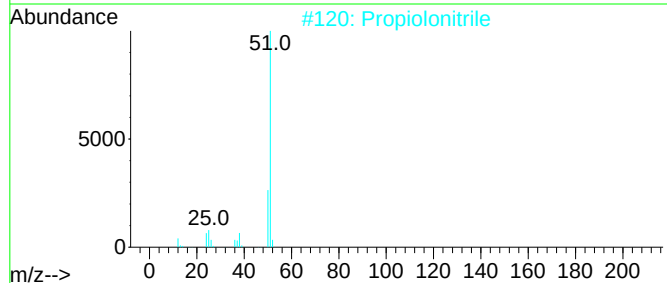
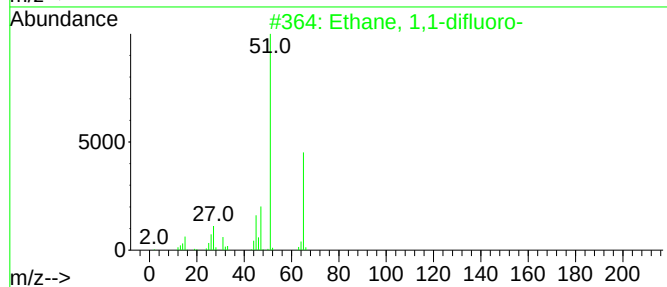
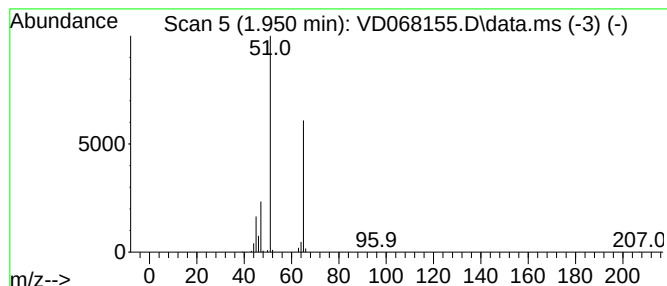
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D011921S.M
Quant Title : SW846 8260

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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.950	17.09 ug/l	558836	Pentafluorobenzene	7.985

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			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-dif...	1.950	17.1	ug/l	558836	1	7.985	1635100	50.0