

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_D\DATA\VD010620\
 Data File : VD064782.D
 Acq On : 06 Jan 2020 19:34
 Operator : VA/SY
 Sample : K6482-05
 Misc : 7.23G/5.00ml/MSVOA_D/SOIL
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 MC-123119-0-1-GRAB-B4

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\82D122419S.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	1.956	3	6	15	rVB	200066	282865	11.19%	2.076%
2	2.344	69	72	78	rVB3	16398	25980	1.03%	0.191%
3	4.967	506	518	528	rBV3	18746	62276	2.46%	0.457%
4	7.920	1010	1020	1025	rBV	351253	849459	33.59%	6.236%
5	7.985	1025	1031	1041	rVB	632926	1397434	55.26%	10.258%
6	8.332	1079	1090	1100	rBV	300454	668065	26.42%	4.904%
7	8.867	1172	1181	1194	rBV	916761	1820593	72.00%	13.364%
8	10.344	1422	1432	1450	rBV	1405588	2528640	100.00%	18.562%
9	11.649	1646	1654	1667	rBV	1234385	2157355	85.32%	15.836%
10	11.855	1681	1689	1699	rBV3	52860	115087	4.55%	0.845%
11	12.185	1732	1745	1752	rBV4	24422	60094	2.38%	0.441%
12	12.638	1814	1822	1830	rBV	1004235	1621425	64.12%	11.902%
13	13.585	1974	1983	1992	rBV	1212133	2033534	80.42%	14.927%

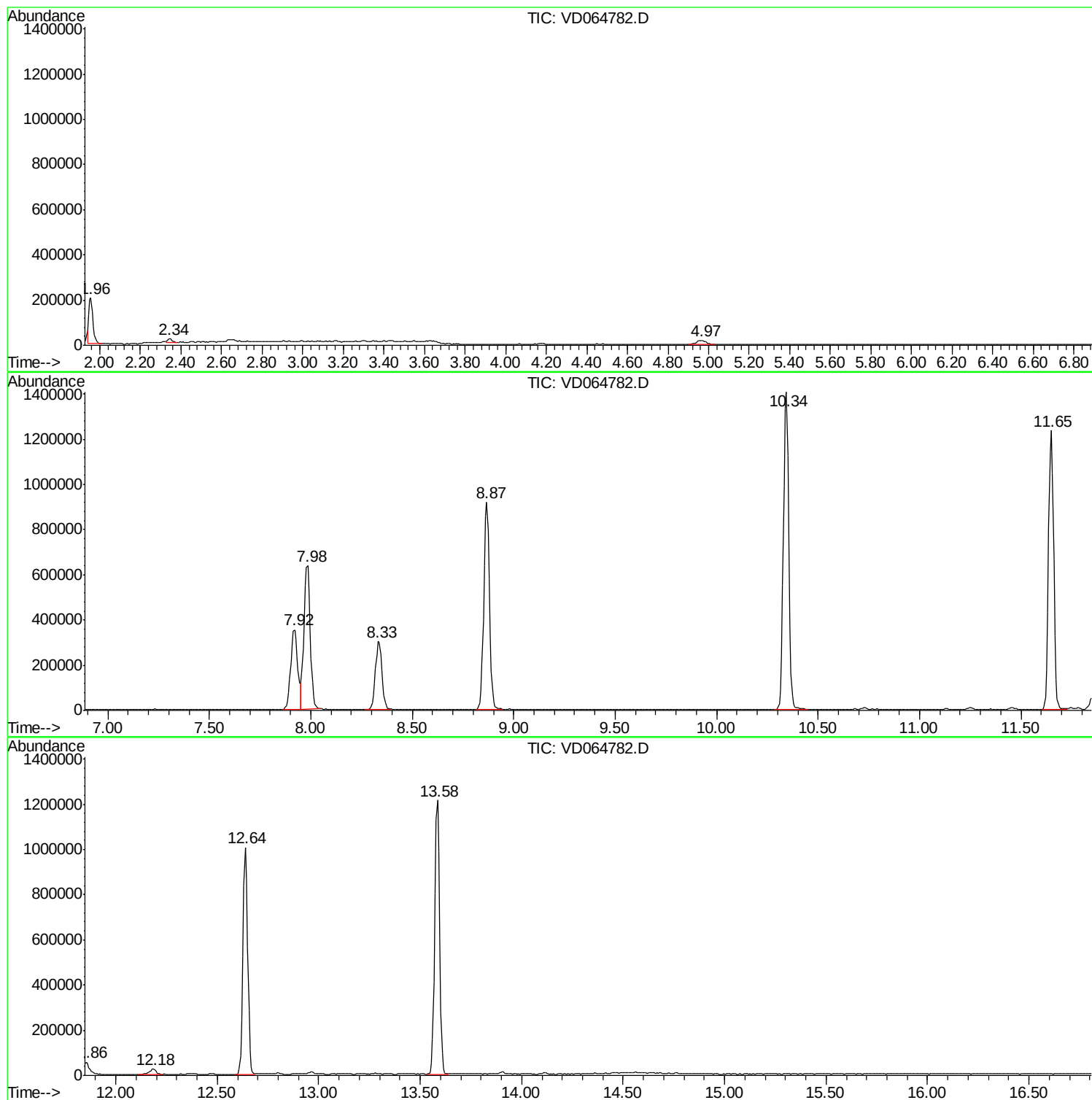
Sum of corrected areas: 13622807

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Quant Method : Z:\V0ASRV\HPCHEM1\MSV0A_D\METHOD\82D122419S.M
Quant Title : SW846 8260

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



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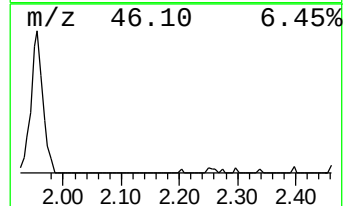
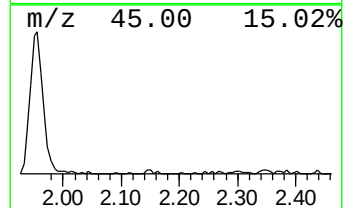
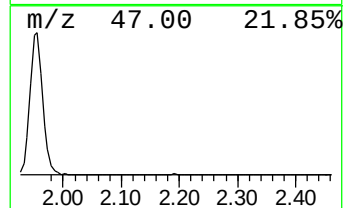
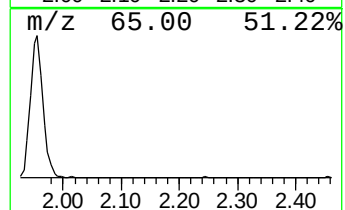
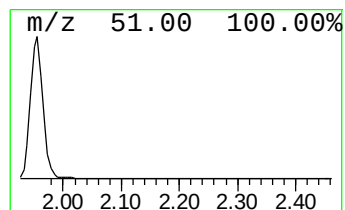
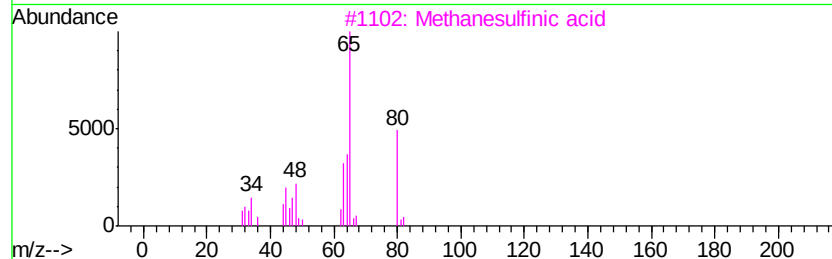
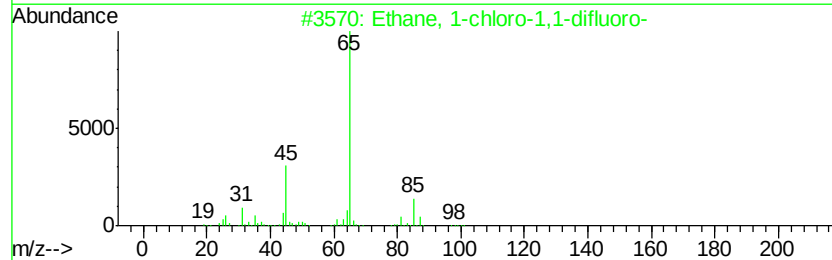
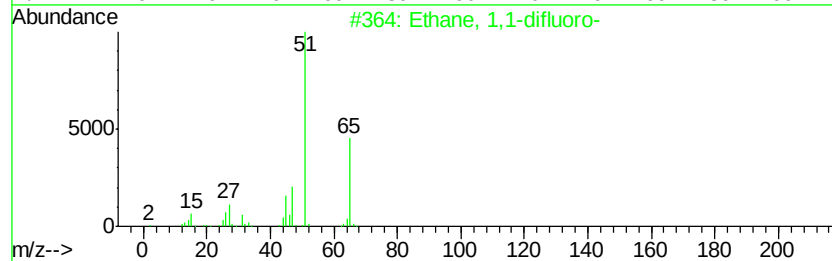
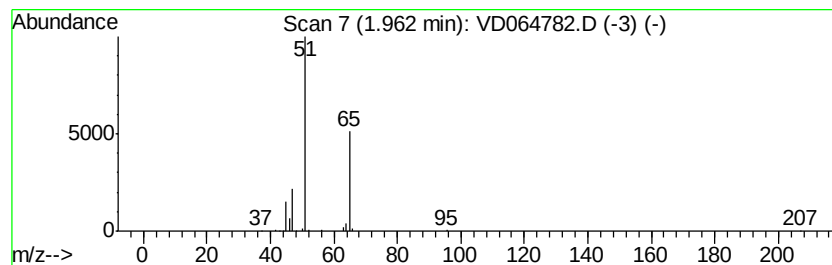
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Quant Title : SW846 8260

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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.
1.96	10.12 ug/l	282865	Pentafluorobenzene	7.98

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Ethane, 1,1-difluoro-	66	C2H4F2	000075-37-6	91
2		Ethane, 1-chloro-1,1-difluoro-	100	C2H3ClF2	000075-68-3	9
3		Methanesulfinic acid	80	CH4O2S	017696-73-0	9
4		2-p-Nitrobenzoyl-1,3,5-tribenzyl...	569	C33H31N08	1000129-85-6	4
5		Ethanamine, N-chloro-N,1,1-trifl...	133	C2H3ClF3N	016276-45-2	4



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Ethane, 1,1-diflu...	1.96	10.1	ug/l	282865	1	7.98	1397430	50.0