

Data Path : Z:\VOASRV\HPCHEM1\MSVOA D\DATA\VD041219\
 Data File : VD061789.D
 Acq On : 12 Apr 2019 20:23
 Operator : VA/AP
 Sample : K2360-02
 Misc : 5.14g/5ml/MSVOA D/SOIL
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 FK-BPM-01-E2-E2A-F2-F2A

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\82D040619S.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	1.491	3	7	11	rVB	178010	493972	23.62%	3.710%
2	1.550	11	13	29	rVB	77387	232042	11.10%	1.743%
3	3.214	177	182	188	rBV	15510	37131	1.78%	0.279%
4	3.804	236	242	251	rBV	46333	133348	6.38%	1.002%
5	4.690	327	332	337	rBV	8767	22805	1.09%	0.171%
6	6.915	550	558	564	rBV	345576	998494	47.75%	7.500%
7	7.033	564	570	591	rVB	509452	1572343	75.20%	11.810%
8	7.437	605	611	623	rBV	217500	576487	27.57%	4.330%
9	8.205	683	689	704	rBV	759888	1899748	90.86%	14.269%
10	10.400	906	912	925	rBV	849985	1889274	90.36%	14.190%
11	12.359	1107	1111	1120	rBV	1261937	2090908	100.00%	15.705%
12	13.550	1228	1232	1242	rVB	955802	1325660	63.40%	9.957%
13	13.708	1244	1248	1251	rBB	21642	32799	1.57%	0.246%
14	14.515	1326	1330	1333	rBV	1439554	2009044	96.08%	15.090%

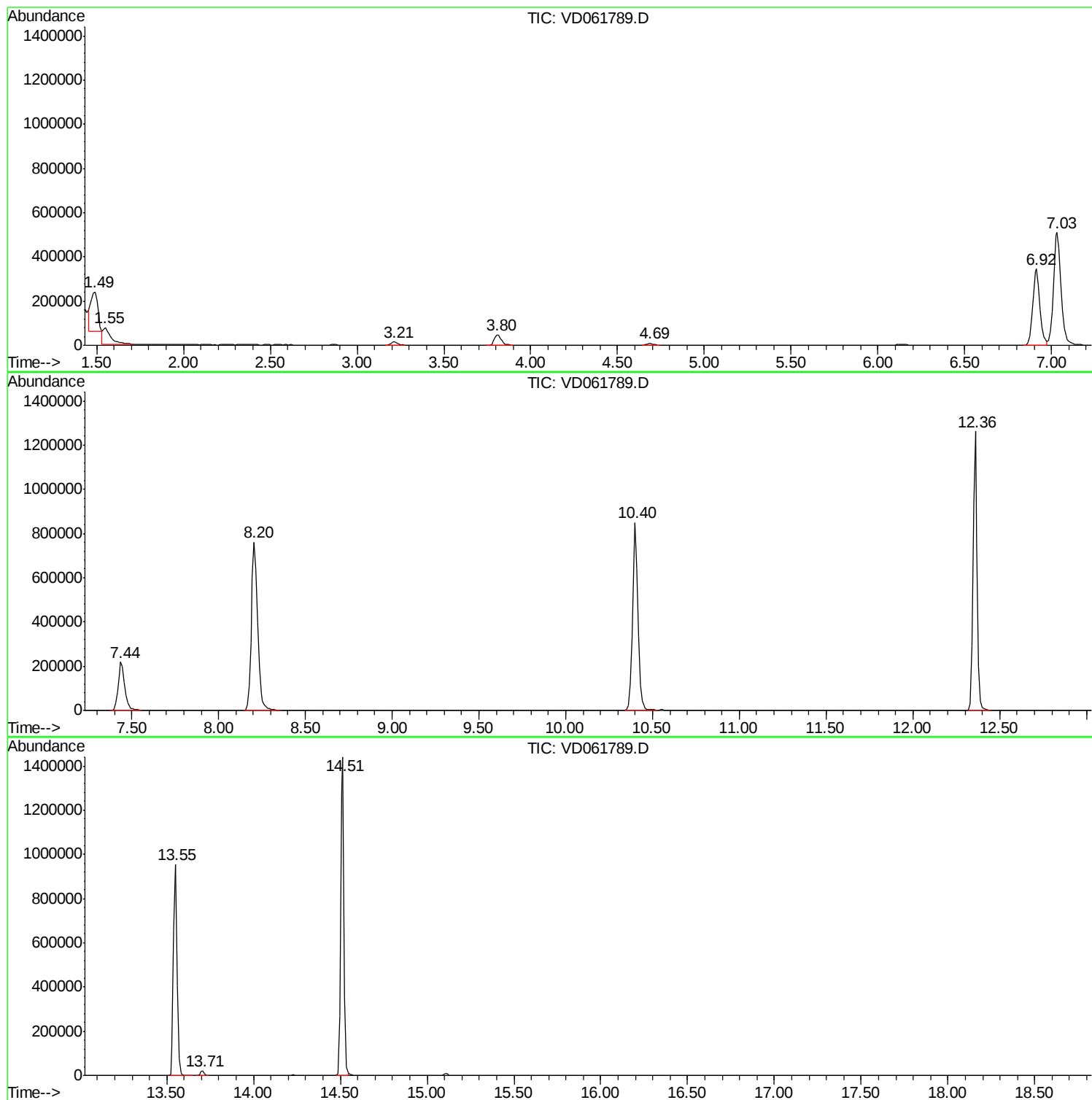
Sum of corrected areas: 13314055

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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D040619S.M
Quant Title : SW846 8260

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



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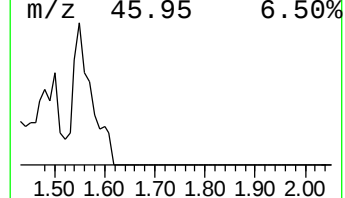
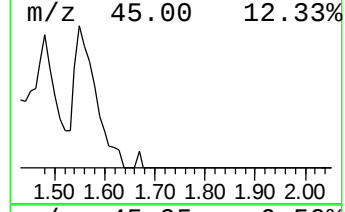
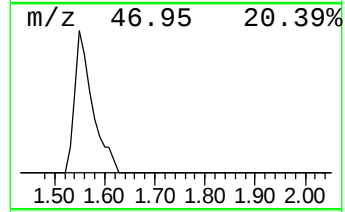
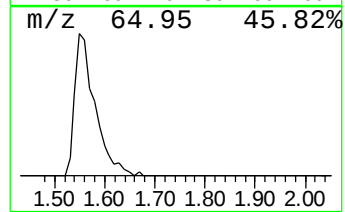
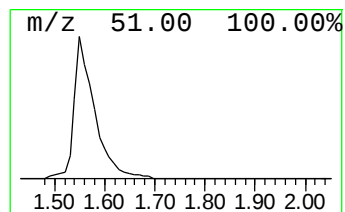
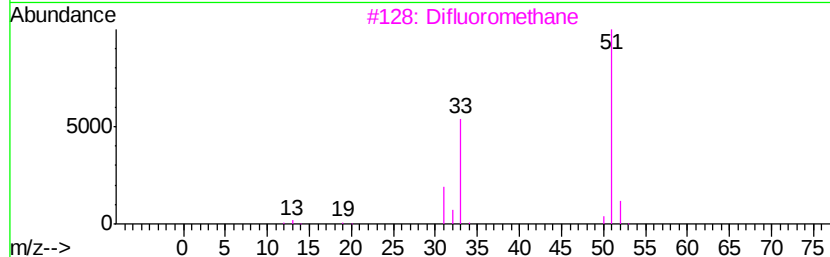
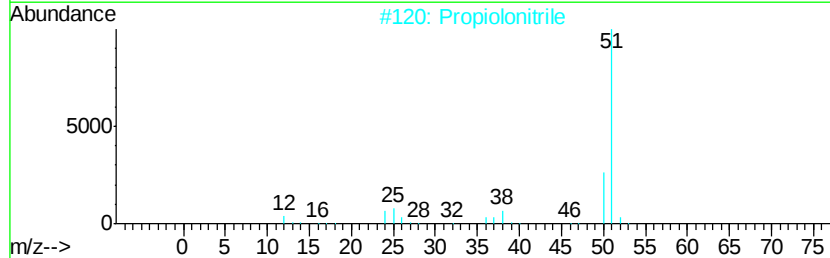
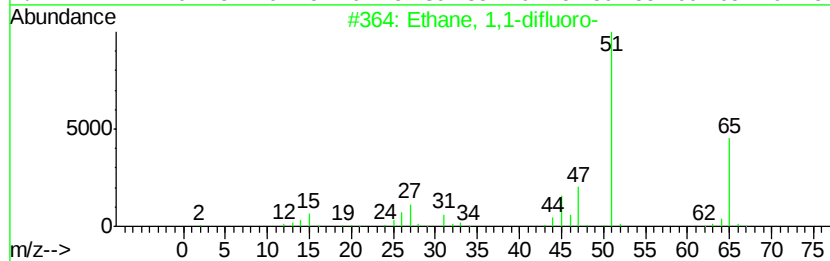
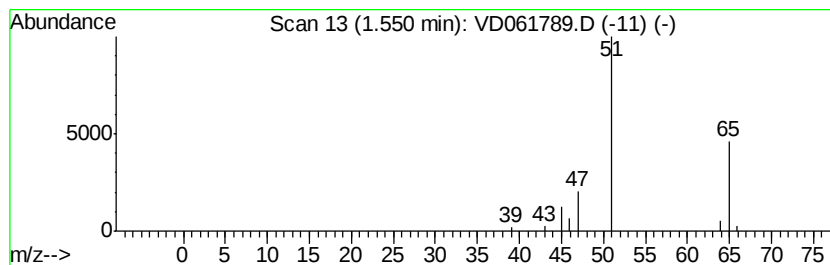
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Quant Title : SW846 8260

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TIC Integration Parameters: LSCINT.P

Peak Number 2 Ethane, 1,1-difluoro- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.55	7.38 ug/l	232042	Pentafluorobenzene	7.03

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethane, 1,1-difluoro-	66	C2H4F2	000075-37-6	90
2			Propiolonitrile	51	C3HN	001070-71-9	3
3			Difluoromethane	52	CH2F2	000075-10-5	1
4			2-Chloroethanol	80	C2H5ClO	000107-07-3	1
5			Propanoyl chloride	92	C3H5ClO	000079-03-8	1



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Ethane, 1,1-diflu...	1.55	7.4	ug/l	232042	1	7.03	1572340	50.0