

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_D\DATA\VD010320\
 Data File : VD064754.D
 Acq On : 04 Jan 2020 02:04
 Operator : VA/SY
 Sample : K6463-01
 Misc : 6.95G/5.00ml/MSVOA_D/SOIL
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 GP-6-(1-3)

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	138832	197237	7.93%	1.521%
2	4.962	506	517	531	rBV4	14166	49249	1.98%	0.380%
3	7.920	1009	1020	1025	rBV	358212	860163	34.57%	6.633%
4	7.979	1025	1030	1041	rVB	648958	1433571	57.61%	11.054%
5	8.338	1079	1091	1101	rBV	307720	689326	27.70%	5.315%
6	8.867	1169	1181	1192	rBV	910407	1824256	73.31%	14.067%
7	10.344	1423	1432	1440	rBV	1378811	2488507	100.00%	19.189%
8	11.649	1646	1654	1667	rBV	1235048	2093328	84.12%	16.142%
9	12.638	1814	1822	1832	rBV	925454	1505149	60.48%	11.606%
10	13.585	1976	1983	1991	rVB	1152757	1827678	73.44%	14.093%

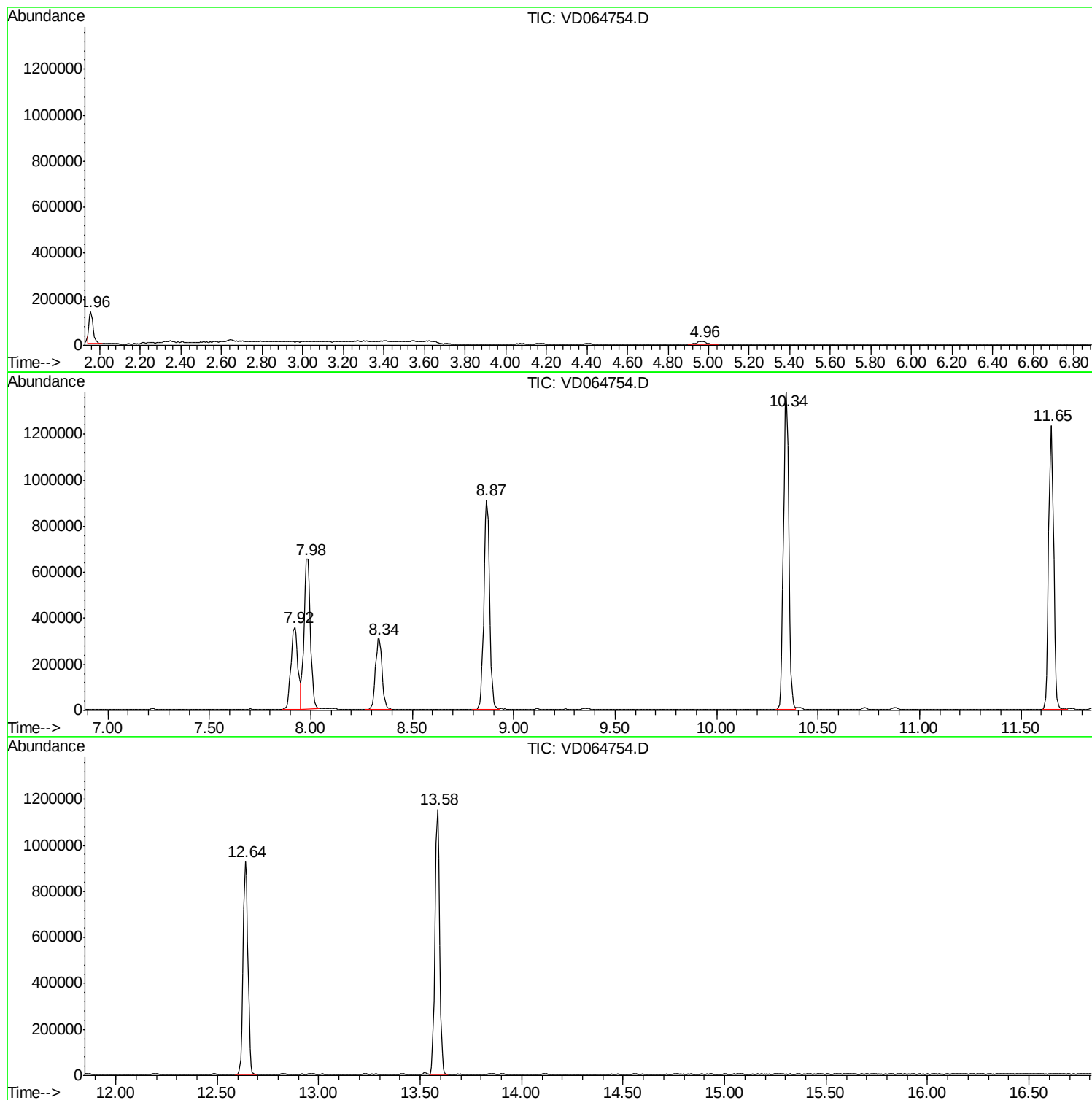
Sum of corrected areas: 12968464

Data Path : Z:\V0ASRV\HPCHEM1\MSV0A_D\DATA\VD010320\
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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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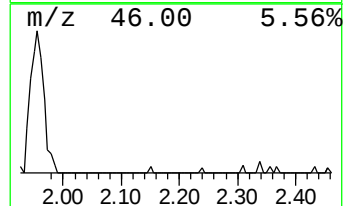
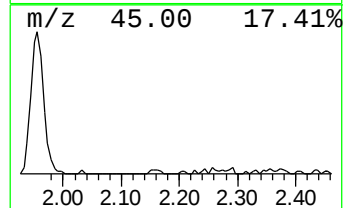
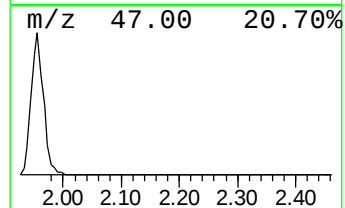
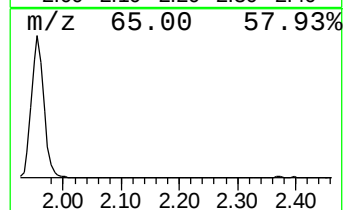
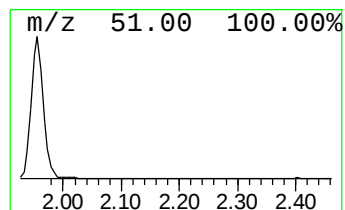
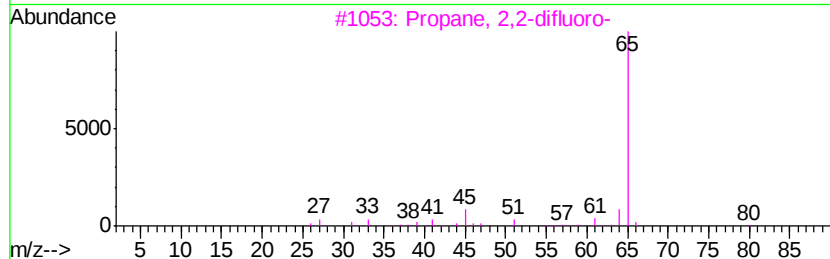
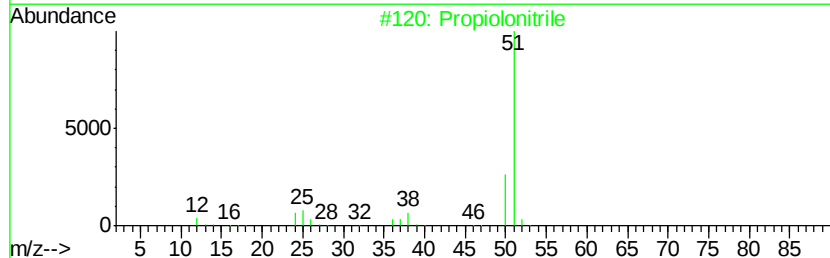
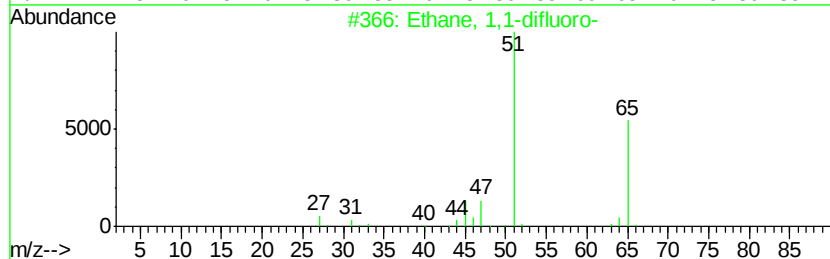
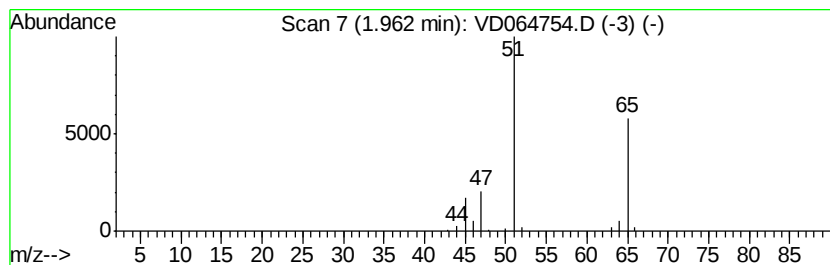
Quant Method : Z:\V0ASRV\HPCHEM1\MSV0A_D\METHOD\82D122419S.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.96	6.88 ug/l	197237	Pentafluorobenzene	7.98

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			Propane, 2,2-difluoro-	80	C3H6F2	000420-45-1	2
4			Difluorochloromethane	86	CHClF2	000075-45-6	2
5			2-Butyne-1,4-diol, diformate	142	C6H6O4	036677-73-3	1



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Ethane, 1,1-diflu...	1.96	6.9	ug/l	197237	1	7.98	1433570	50.0