

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_D\DATA\VD010320\
 Data File : VD064755.D
 Acq On : 04 Jan 2020 02:33
 Operator : VA/SY
 Sample : K6463-02
 Misc : 8.07G/5.00ml/MSVOA_D/SOIL
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 GP-6-(4-6)

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	17	rVB	145409	213381	8.57%	1.654%
2	4.961	508	517	527	rBV3	16837	50591	2.03%	0.392%
3	7.920	1011	1020	1025	rBV	343880	829131	33.30%	6.428%
4	7.985	1025	1031	1041	rVB	640628	1418609	56.97%	10.998%
5	8.338	1081	1091	1101	rBV	292483	639477	25.68%	4.958%
6	8.867	1173	1181	1194	rBV	902820	1780407	71.50%	13.803%
7	10.343	1422	1432	1447	rBV	1395724	2490190	100.00%	19.306%
8	11.649	1646	1654	1666	rBV	1224998	2078374	83.46%	16.113%
9	12.637	1812	1822	1833	rVB	924812	1522262	61.13%	11.802%
10	13.584	1976	1983	1991	rVB	1172549	1876407	75.35%	14.547%

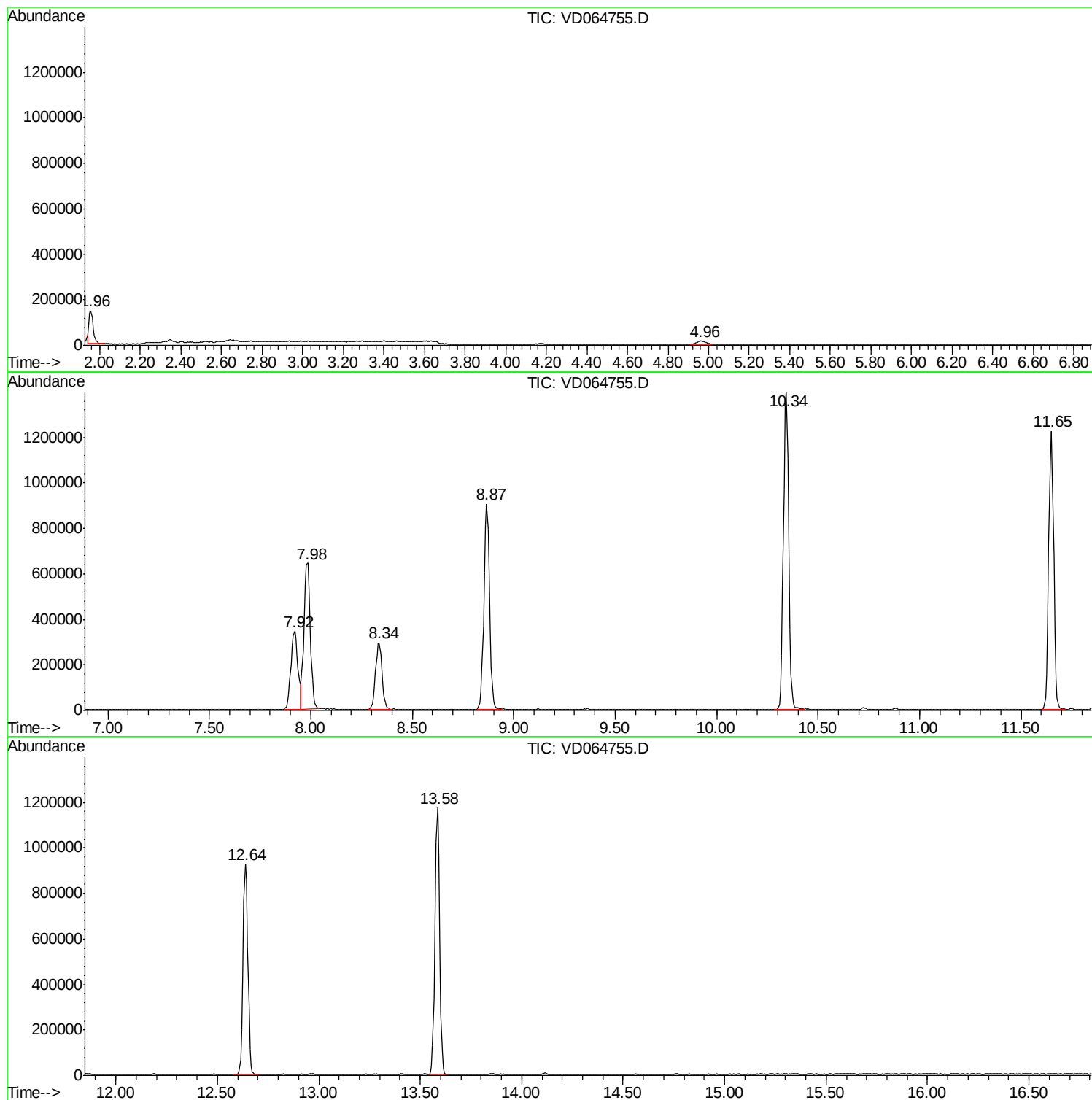
Sum of corrected areas: 12898829

Data Path : Z:\V0ASRV\HPCHEM1\MSV0A_D\DATA\VD010320\
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Instrument :
MSV0A_D
ClientSampleId :
GP-6-(4-6)

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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Instrument :
MSV0A_D
ClientSampled :
GP-6-(4-6)

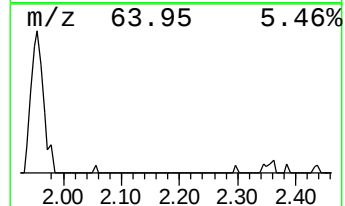
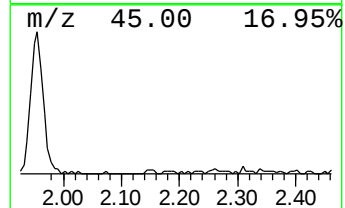
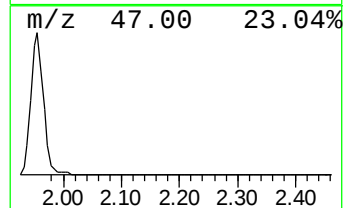
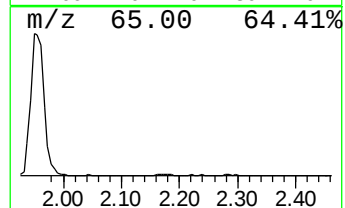
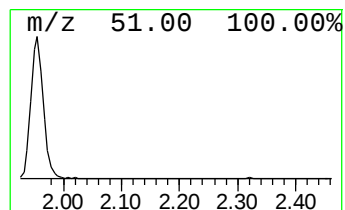
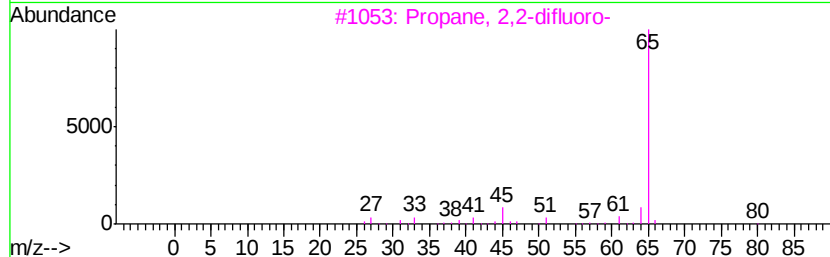
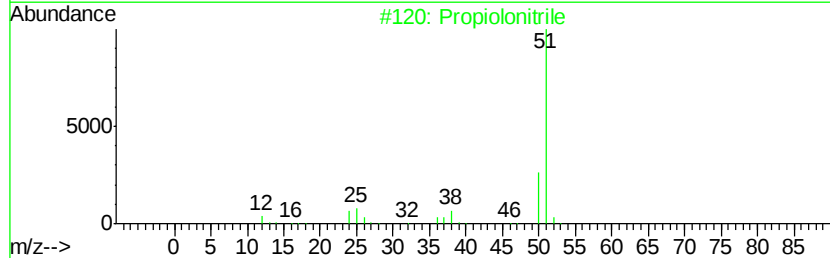
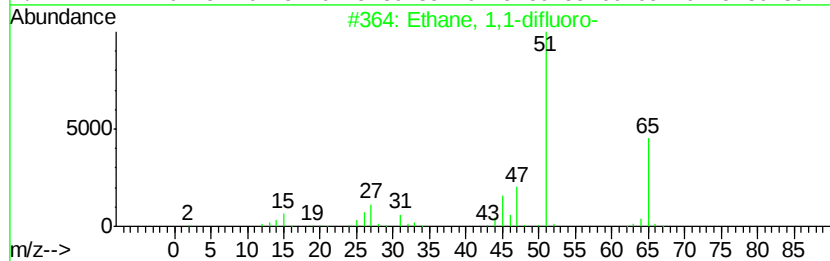
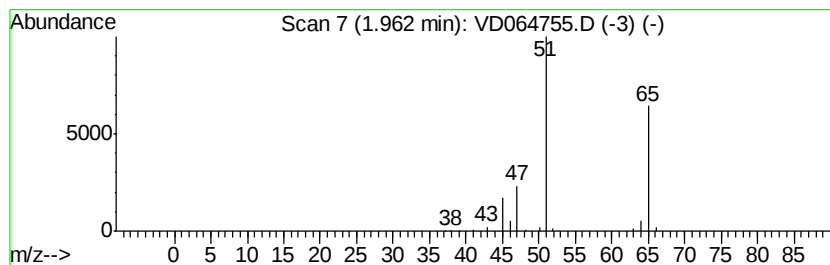
Quant Method : Z:\V0ASRV\HPCHEM1\MSV0A_D\METHOD\82D122419S.M
Quant Title : SW846 8260

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	7.52 ug/l	213381	Pentafluorobenzene	7.98

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...	1.96	7.5	ug/l	213381	1	7.98	1418610	50.0