

Data Path : Z:\VOASRV\HPCHEM1\MSVOA D\DATA\VD021320\
 Data File : VD065179.D
 Acq On : 13 Feb 2020 16:17
 Operator : VA/SY
 Sample : L1529-04
 Misc : 4.47G/5.00ml/MSVOA D/SOIL
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 FK-GF-02-049-B

Integration Parameters: RTEINT.P

Integrator: RTE

Smoothing : ON

Sampling : 1

Start Thrs: 0.2

Stop Thrs : 0

Filtering: 5

Min Area: 3 % of largest Peak

Max Peaks: 100

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\VOASRV\HPCHEM1\MSVOA_D\METHOD\82D020520S.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	1.950	3	5	16	rVB	31353	48182	5.52%	1.009%
2	3.656	293	295	304	rVB5	5850	9832	1.13%	0.206%
3	4.162	373	381	391	rBV3	12349	34485	3.95%	0.722%
4	7.215	893	900	908	rVB	10415	23478	2.69%	0.492%
5	7.920	1009	1020	1025	rBV	126146	295304	33.81%	6.186%
6	7.985	1025	1031	1041	rVB	185194	420747	48.17%	8.814%
7	8.338	1083	1091	1103	rBV2	121736	276618	31.67%	5.795%
8	8.867	1173	1181	1193	rBV	269486	548647	62.81%	11.494%
9	10.344	1424	1432	1440	rBV	442450	772221	88.40%	16.177%
10	10.726	1491	1497	1504	rBV3	7793	12140	1.39%	0.254%
11	11.650	1646	1654	1669	rVB	487447	831470	95.19%	17.418%
12	12.638	1814	1822	1830	rBV	373190	626890	71.77%	13.133%
13	13.585	1976	1983	1990	rBV	539877	873510	100.00%	18.299%

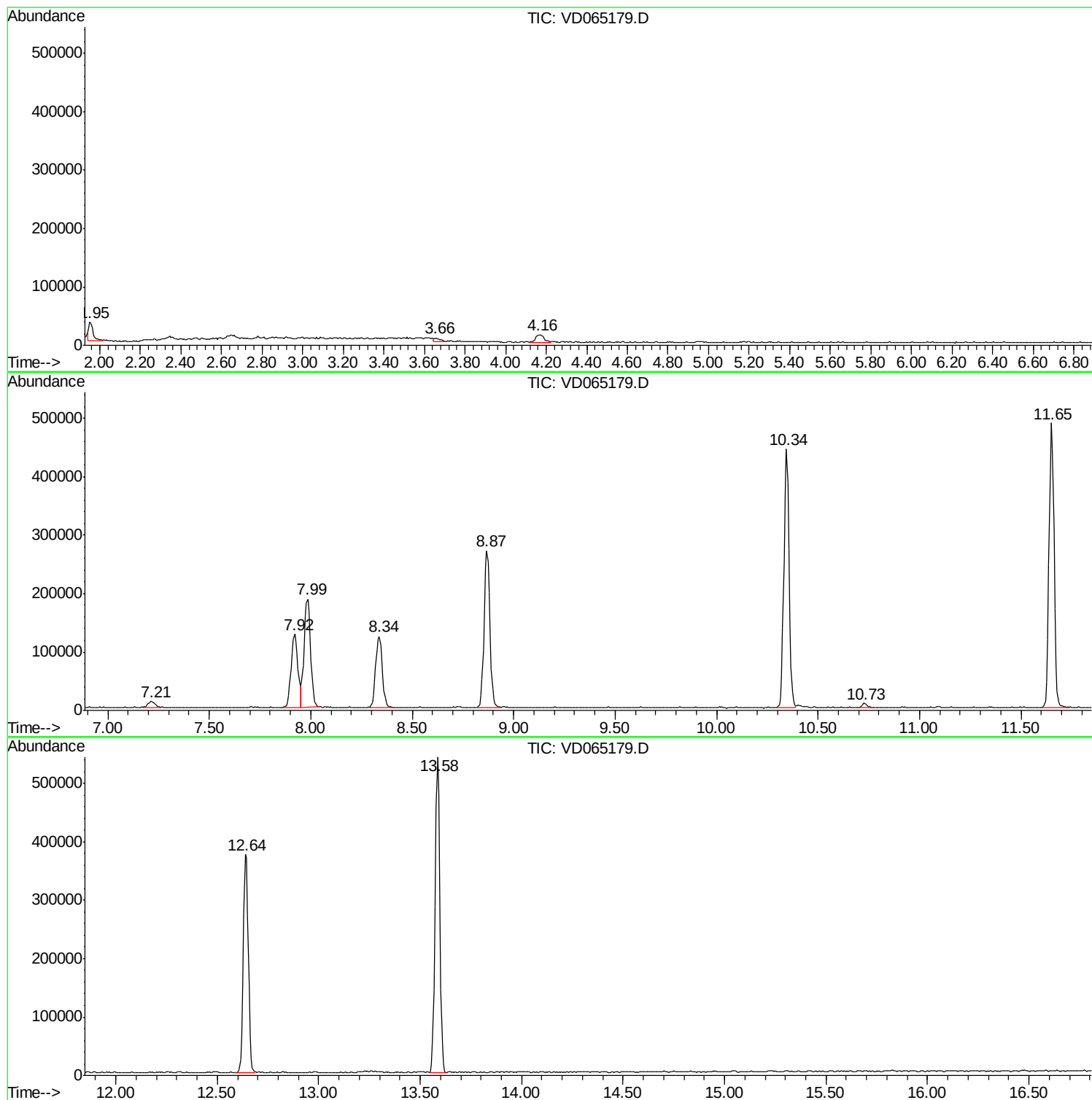
Sum of corrected areas: 4773524

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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D020520S.M
Quant Title : SW846 8260

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



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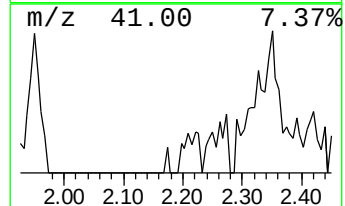
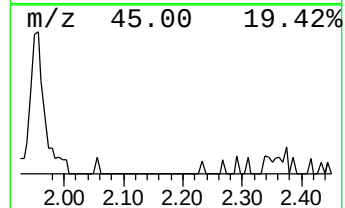
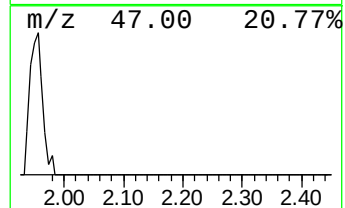
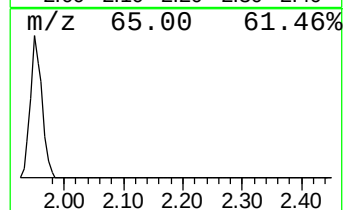
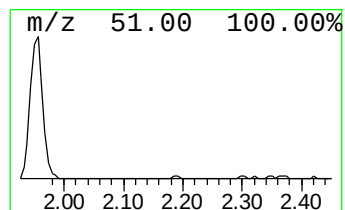
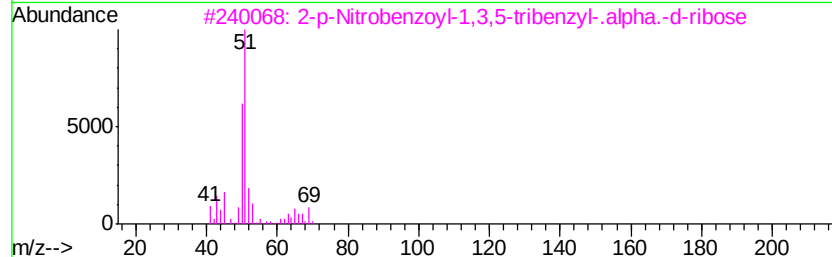
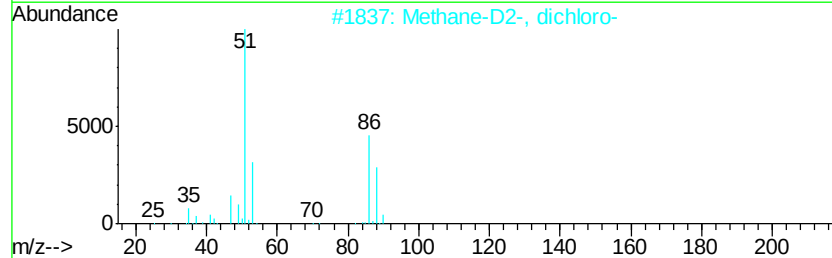
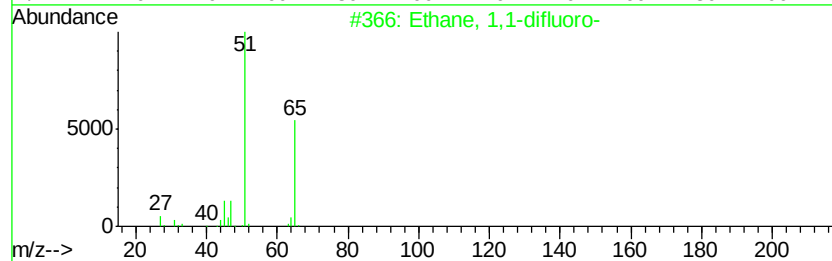
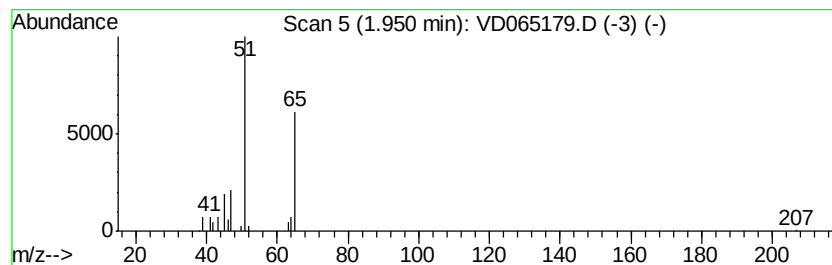
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D020520S.M
Quant Title : SW846 8260

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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.
1.95	5.73 ug/l	48182	Pentafluorobenzene	7.99

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Ethane, 1,1-difluoro-	66	C2H4F2	000075-37-6	74
2		Methane-D2-, dichloro-	86	CD2Cl2	001665-00-5	38
3		2-p-Nitrobenzoyl-1,3,5-tribenzyl...	569	C33H31N08	1000129-85-6	9
4		Ethane, 1-chloro-1,1-difluoro-	100	C2H3ClF2	000075-68-3	4
5		s-Triazolo[4,3-a]pyridine, 7-met...	133	C7H7N3	004919-10-2	4



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TIC Top Hit name	RT	EstConc	Units	Response	---Internal Standard---			
					#	RT	Resp	Conc
Ethane, 1,1-diflu...	1.95	5.7	ug/l	48182	1	7.99	420747	50.0