

Data Path : Z:\VOASRV\HPCHEM1\MSVOA D\DATA\VD040119\
 Data File : VD061479.D
 Acq On : 1 Apr 2019 20:09
 Operator : VA/AP
 Sample : K2182-11
 Misc : 5.30g/5ml/MSVOA D/SOIL
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 S8A(10)

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\82D032719S.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	1.554	11	13	27	rVB	134178	335410	14.49%	2.313%
2	1.850	38	43	50	rVB3	36376	97480	4.21%	0.672%
3	2.667	120	126	132	rVB5	9923	32755	1.42%	0.226%
4	3.228	178	183	190	rBV	11961	33746	1.46%	0.233%
5	3.828	237	244	259	rBV	72371	227305	9.82%	1.568%
6	6.919	551	558	565	rBV	393188	1156163	49.95%	7.975%
7	7.047	565	571	588	rVB	635446	1859569	80.34%	12.826%
8	7.451	606	612	626	rBV	265302	704521	30.44%	4.859%
9	8.219	684	690	706	rBV	864478	2218395	95.84%	15.301%
10	10.404	906	912	919	rBV	925859	2103311	90.87%	14.507%
11	10.503	919	922	927	rVB	74717	151136	6.53%	1.042%
12	12.363	1107	1111	1124	rBV	1357631	2314755	100.00%	15.966%
13	13.555	1228	1232	1237	rBV	995456	1362913	58.88%	9.401%
14	13.712	1245	1248	1251	rBV	27154	38405	1.66%	0.265%
15	14.519	1326	1330	1333	rBV	1453092	1862360	80.46%	12.845%

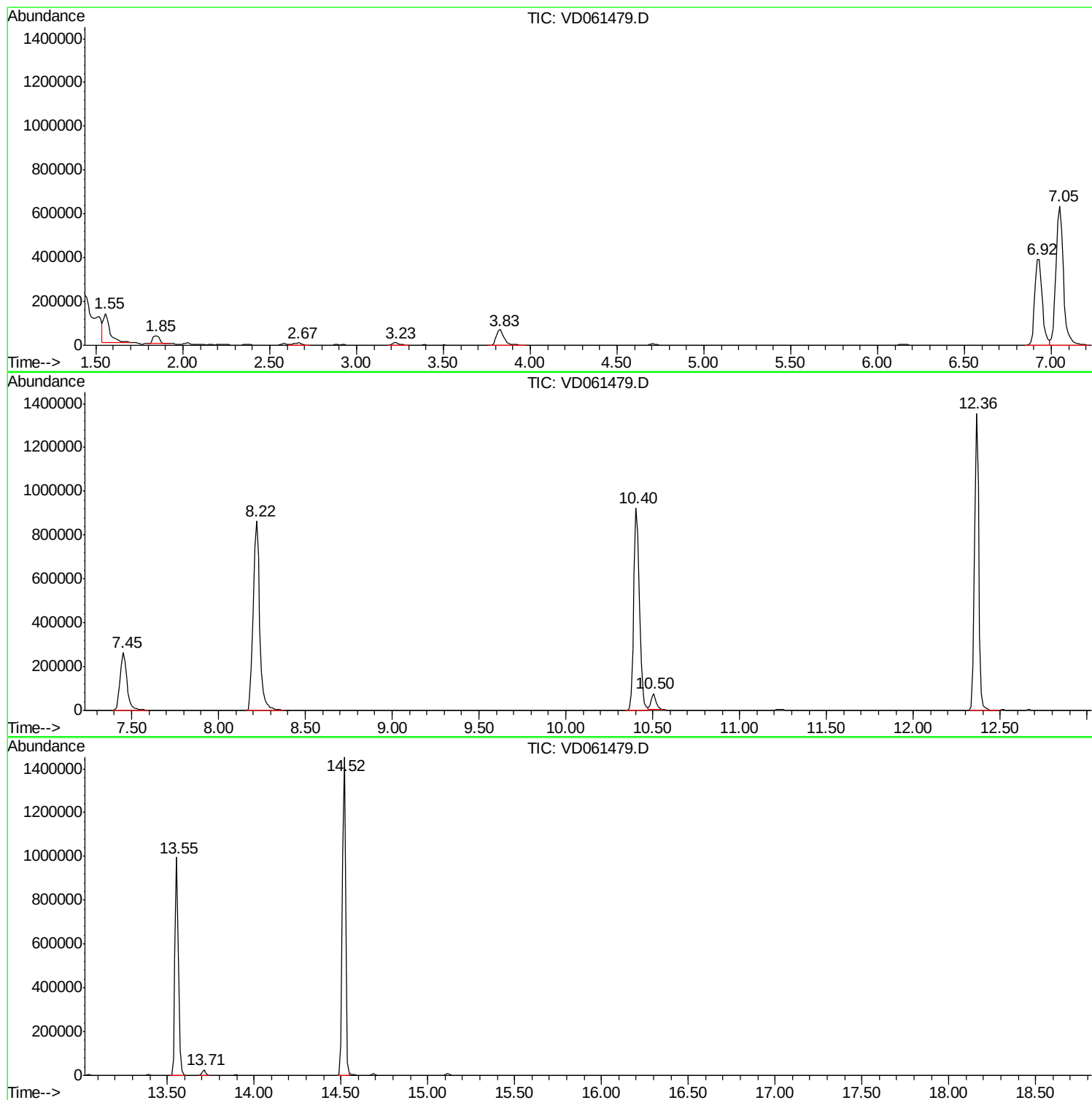
Sum of corrected areas: 14498224

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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D032719S.M
Quant Title : SW846 8260

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



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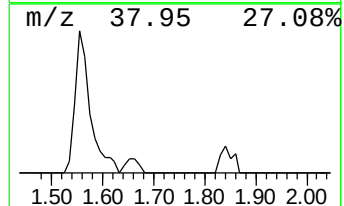
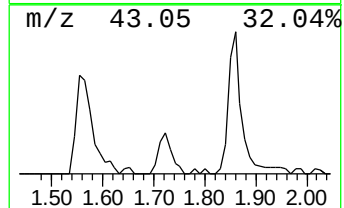
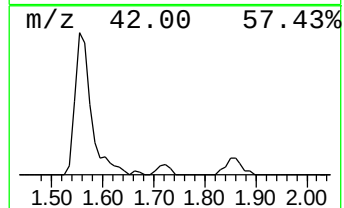
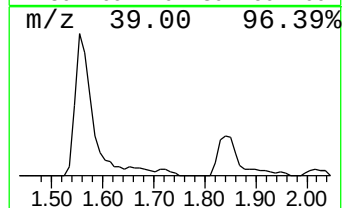
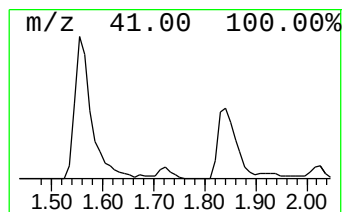
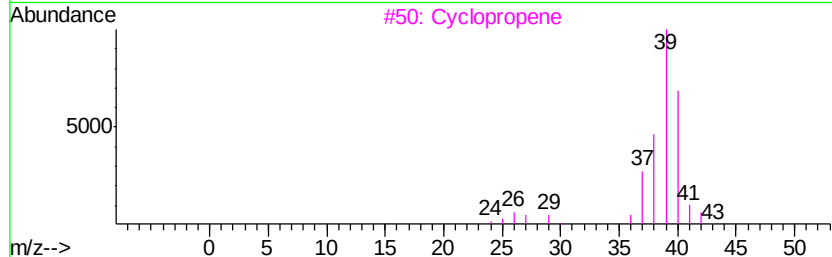
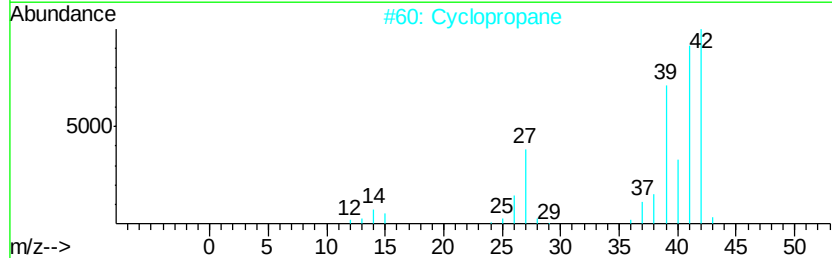
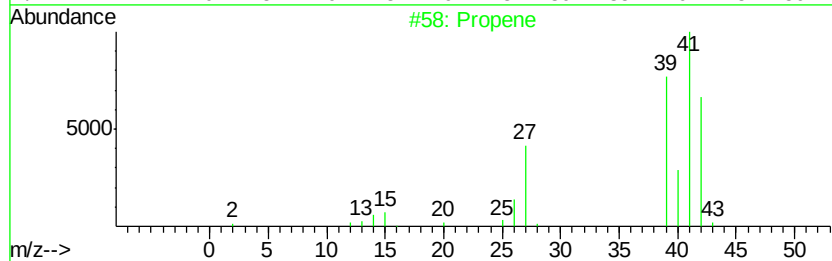
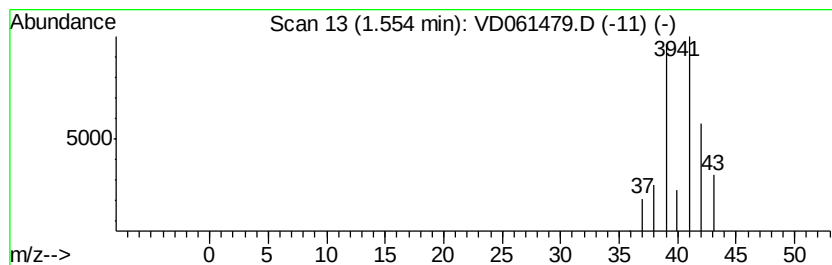
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Peak Number 1 unknown1.55 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.	
1.55	9.02 ug/l	335410	Pentafluorobenzene	7.05	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Propene	42	C3H6	000115-07-1	42
2	Cyclopropane	42	C3H6	000075-19-4	4
3	Cyclopropene	40	C3H4	002781-85-3	4
4	2-Oxetanone, 4-methyl-	86	C4H6O2	003068-88-0	1
5	Butanenitrile	69	C4H7N	000109-74-0	1



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown1.55	1.55	9.0	ug/l	335410	1	7.05	1859570	50.0