

Data Path : Z:\VOASRV\HPCHEM1\MSVOA W\DATA\VW082019\
 Data File : VW012041.D
 Acq On : 20 Aug 2019 20:30
 Operator : SY/VA
 Sample : K4387-04
 Misc : 6.46G/5ML/MSVOA W/SOIL
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 T6-C-(0)

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_W\METHOD\82W072719S.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	4.916	481	494	504	rBV3	17571	63592	4.78%	0.958%
2	7.543	916	925	936	rBV4	6480	18720	1.41%	0.282%
3	7.885	969	981	986	rBV2	172959	413405	31.08%	6.226%
4	7.946	986	991	1002	rVB	330574	712164	53.54%	10.725%
5	8.305	1039	1050	1060	rBV	187549	417955	31.42%	6.294%
6	8.842	1127	1138	1150	rBV	497612	998499	75.07%	15.037%
7	10.323	1372	1381	1389	rBV	750000	1330053	100.00%	20.030%
8	10.701	1436	1443	1450	rBV2	7520	13764	1.03%	0.207%
9	11.628	1587	1595	1607	rBV	588524	1024721	77.04%	15.432%
10	12.615	1750	1757	1768	rBV	430634	724975	54.51%	10.918%
11	13.554	1905	1911	1923	rBV	417868	706256	53.10%	10.636%
12	14.072	1992	1996	2002	rVB2	16913	26738	2.01%	0.403%
13	14.725	2099	2103	2110	rBV3	16557	24978	1.88%	0.376%
14	15.816	2276	2282	2289	rBV2	84958	164563	12.37%	2.478%

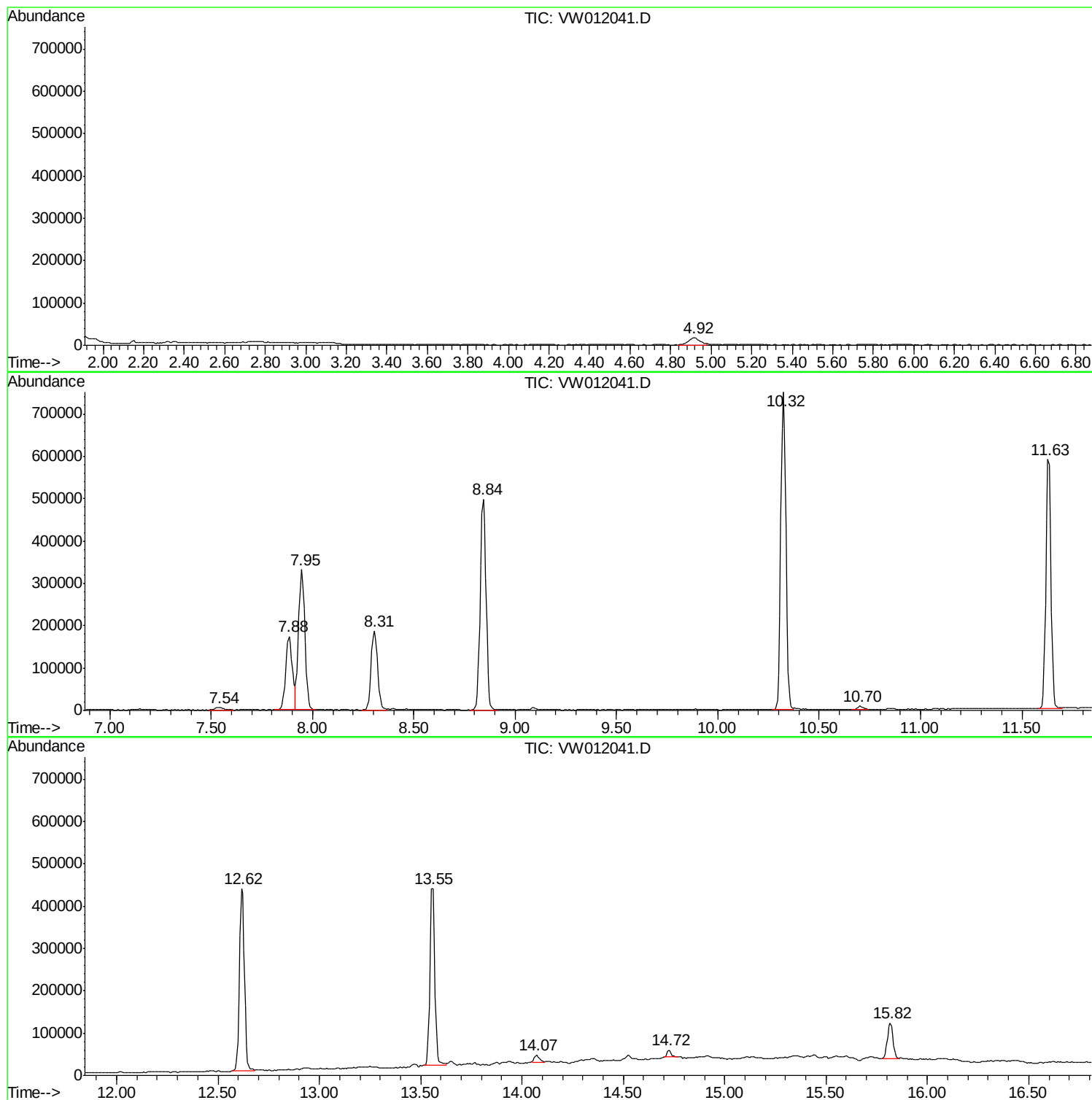
Sum of corrected areas: 6640383

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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W072719S.M
Quant Title : SW846 8260

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



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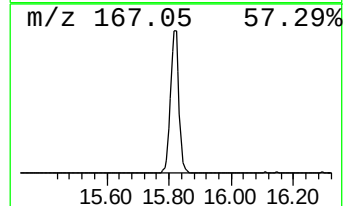
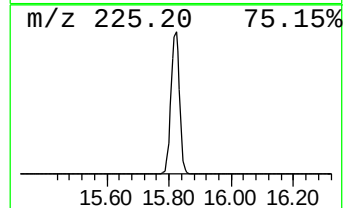
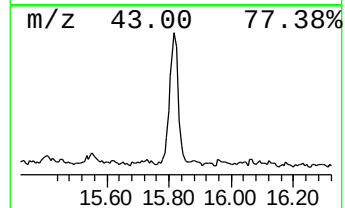
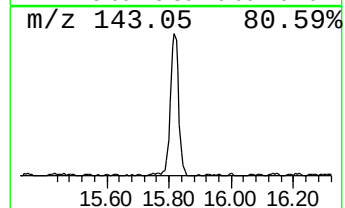
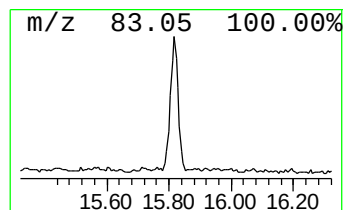
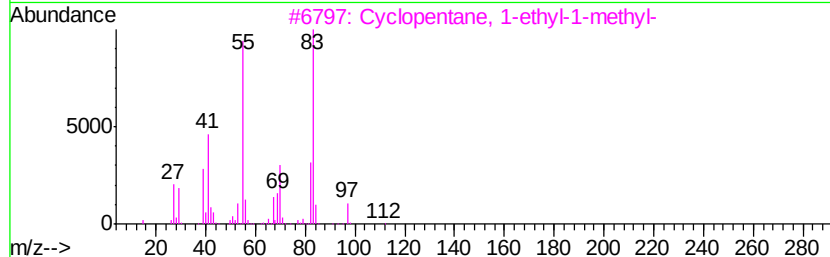
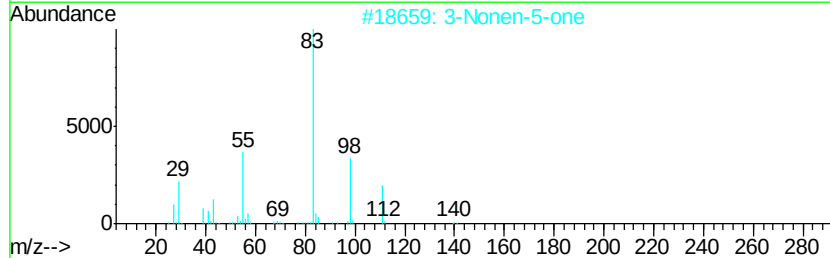
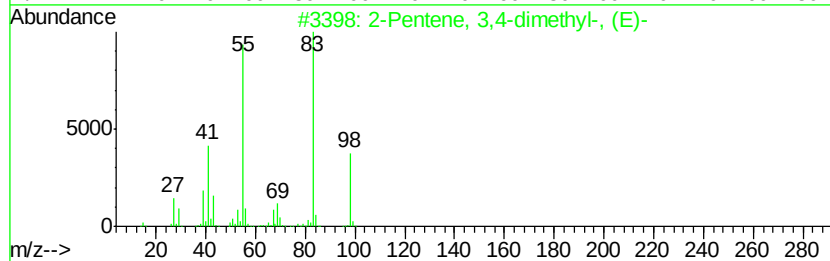
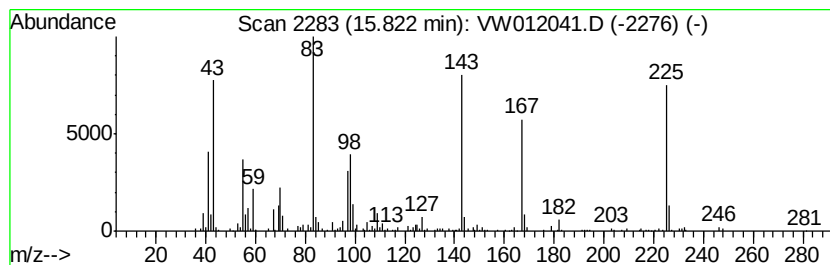
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Peak Number 1 unknown15.82 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.82	11.65 ug/l	164563	1,4-Dichlorobenzene-d4	13.56

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentene, 3,4-dimethyl-, (E)-	98	C7H14	004914-92-5	25
2		3-Nonen-5-one	140	C9H16O	082456-34-6	25
3		Cyclopentane, 1-ethyl-1-methyl-	112	C8H16	016747-50-5	25
4		2-Pentene, 4,4-dimethyl-, (E)-	98	C7H14	000690-08-4	25
5		2-Pentene, 4,4-dimethyl-, (Z)-	98	C7H14	000762-63-0	25



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown15.82	15.82	11.7	ug/l	164563	4	13.56	706256	50.0