

Data Path : Z:\VOASRV\HPCHEM1\MSVOA W\DATA\VW081719\
 Data File : VW011964.D
 Acq On : 16 Aug 2019 14:52
 Operator : SY/VA
 Sample : K4366-01
 Misc : 5.74G/5ML/MSVOA W/SOIL
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 PX-1

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	2.154	35	41	48	rBV	25448	54229	3.86%	0.615%
2	2.660	114	124	132	rBV4	12629	39365	2.80%	0.446%
3	4.909	481	493	509	rVB3	9173	34435	2.45%	0.390%
4	7.884	972	981	986	rBV	166802	401370	28.54%	4.550%
5	7.945	986	991	1003	rVB	314534	682079	48.50%	7.732%
6	8.305	1040	1050	1064	rBV	201976	459612	32.68%	5.210%
7	8.842	1129	1138	1145	rBV	494903	977865	69.53%	11.085%
8	10.323	1374	1381	1389	rBV	769219	1352879	96.20%	15.336%
9	10.707	1440	1444	1451	rVB	16601	28134	2.00%	0.319%
10	11.634	1589	1596	1604	rVB	634039	1095077	77.87%	12.414%
11	12.621	1750	1758	1765	rBV	579836	1229452	87.42%	13.937%
12	13.560	1905	1912	1932	rVV	644333	1406300	100.00%	15.942%
13	13.761	1935	1945	1960	rVB3	191643	797033	56.68%	9.035%
14	13.932	1968	1973	1974	rBV4	38838	66646	4.74%	0.755%
15	15.724	2260	2267	2277	rVB4	59857	197032	14.01%	2.234%

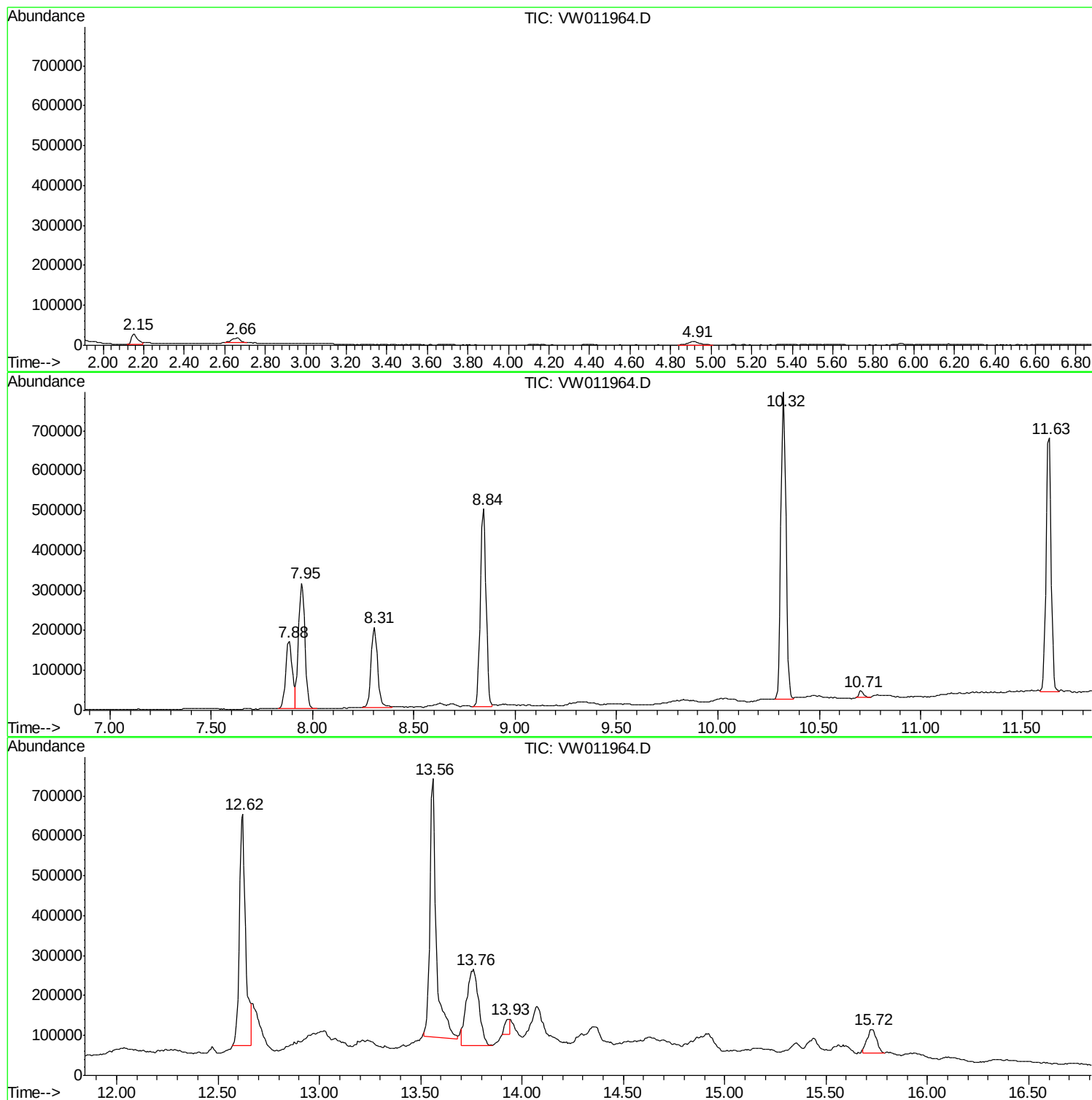
Sum of corrected areas: 8821508

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



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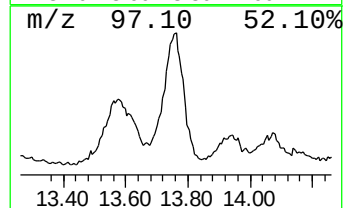
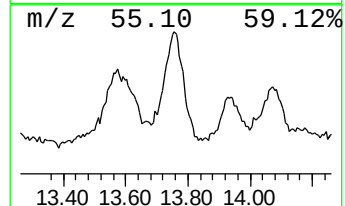
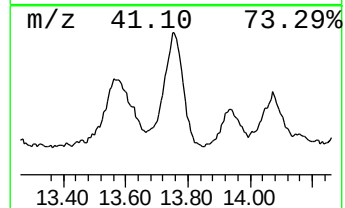
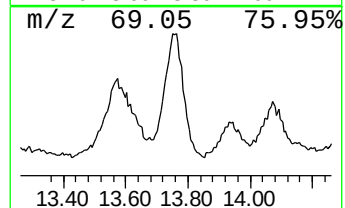
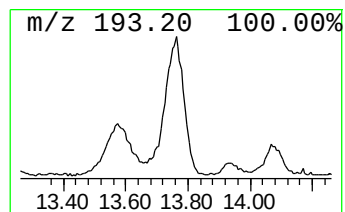
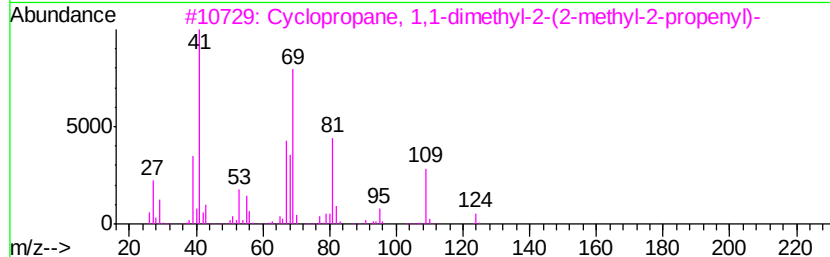
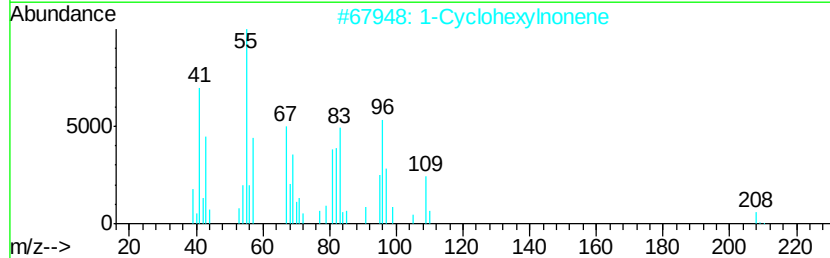
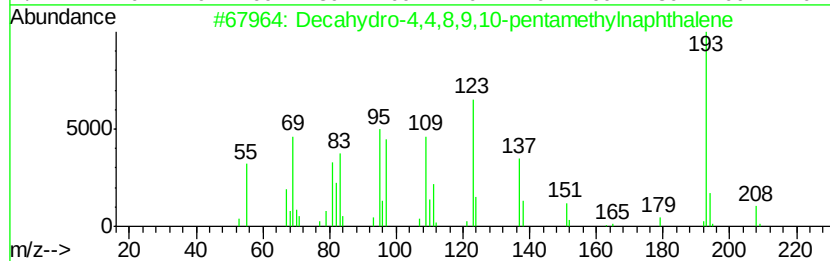
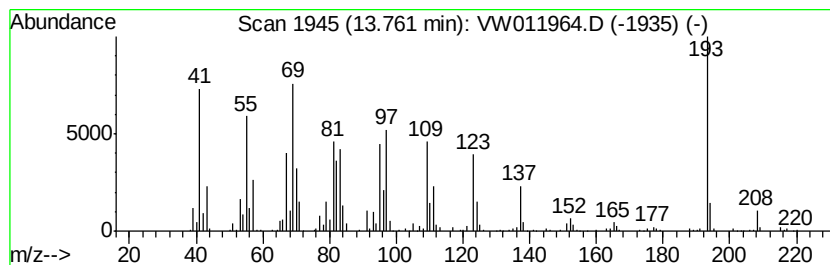
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TIC Integration Parameters: LSCINT.P

Peak Number 1 Decahydro-4,4,8,9,10-pentam... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.		
13.76	28.34 ug/l	797033	1,4-Dichlorobenzene-d4	13.56		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Decahydro-4,4,8,9,10-pentamethyl...	208	C15H28	080655-44-3	76
2		1-Cyclohexylnonene	208	C15H28	114614-84-5	25
3		Cyclopropane, 1,1-dimethyl-2-(2-...	124	C9H16	069147-03-1	22
4		1-(p-Fluorophenyl)-4-piperidone	193	C11H12FNO	1000238-56-7	22
5		Cyclopentanecarboxylic acid, 3-p...	324	C21H40O2	1000280-59-1	14



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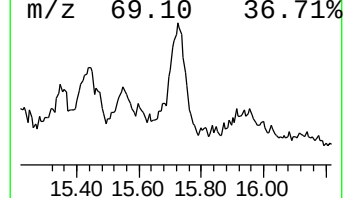
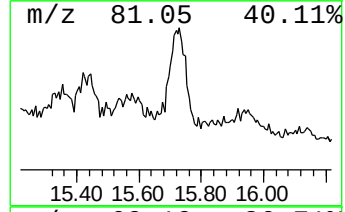
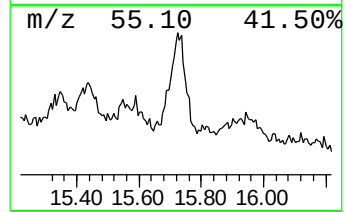
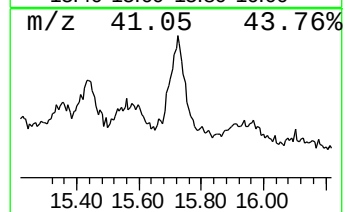
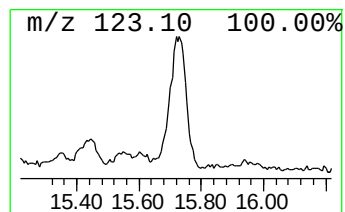
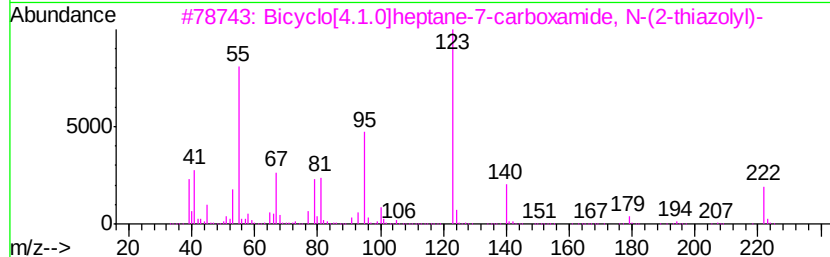
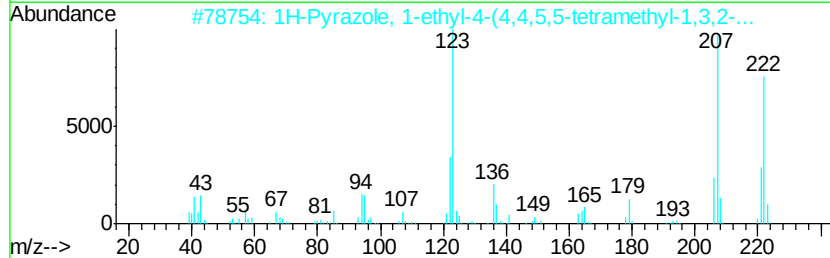
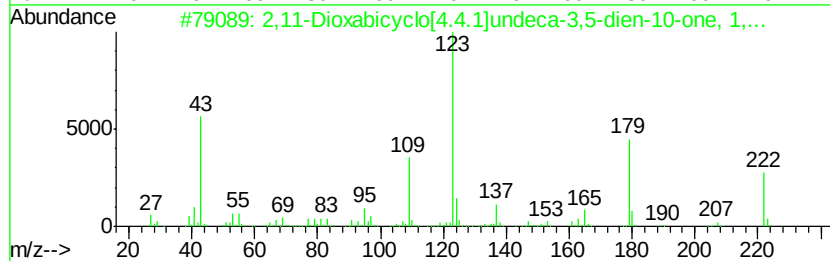
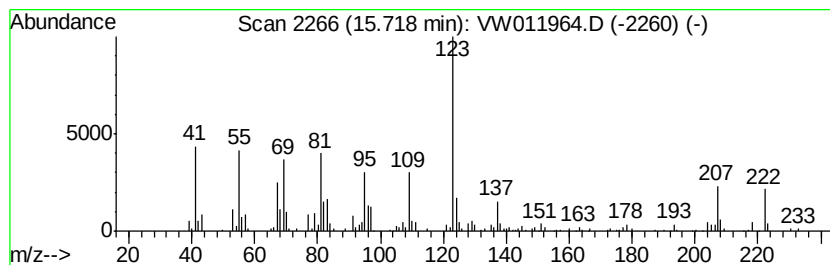
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Peak Number 2 2,11-Dioxabicyclo[4.4.1]und... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.		
15.72	7.01 ug/l	197032	1,4-Dichlorobenzene-d4	13.56		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,11-Dioxabicyclo[4.4.1]undeca-3...	222	C13H18O3	070412-52-1	58	
2	1H-Pyrazole, 1-ethyl-4-(4,4,5,5-...	222	C11H19BN2O2	1000319-69-6	52	
3	Bicyclo[4.1.0]heptane-7-carboxam...	222	C11H14N2OS	329912-72-3	49	
4	3-Fluorobenzoic acid, propargyl ...	178	C10H7FO2	1000325-53-9	43	
5	Ethanone, 1-[3-[2-methyl-2-(5-me...	222	C13H18O3	080114-28-9	43	



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Decahydro-4,4,8,9...	13.76	28.3	ug/l	797033	4	13.56	1406300	50.0
2,11-Dioxabicyclo...	15.72	7.0	ug/l	197032	4	13.56	1406300	50.0