

Data Path : Z:\VOASRV\HPCHEM1\MSVOA W\DATA\VW082619\
 Data File : VW012185.D
 Acq On : 26 Aug 2019 21:40
 Operator : SY/VA
 Sample : K4472-02
 Misc : 5.28G/5ML/MSVOA W/SOIL
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 T1-S1-(2.5)

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\82W082319S.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	2.148	35	40	48	rBV	10678	24930	1.75%	0.285%
2	4.909	480	493	507	rBV2	23288	86246	6.06%	0.984%
3	7.543	916	925	932	rBV3	6890	19702	1.38%	0.225%
4	7.878	968	980	985	rBV	154063	366478	25.73%	4.183%
5	7.945	985	991	1002	rVB	294191	660522	46.37%	7.539%
6	8.305	1039	1050	1062	rBV	172367	387001	27.17%	4.417%
7	8.842	1129	1138	1147	rBV	453786	912377	64.06%	10.414%
8	10.323	1373	1381	1387	rBV	703750	1253384	88.00%	14.306%
9	10.384	1387	1391	1398	rVB	57137	104385	7.33%	1.191%
10	11.628	1589	1595	1603	rVB	566805	954097	66.99%	10.890%
11	12.615	1749	1757	1781	rVB2	526865	1424319	100.00%	16.257%
12	13.554	1905	1911	1929	rVB	522338	1164394	81.75%	13.290%
13	13.749	1935	1943	1955	rVB3	249472	986304	69.25%	11.257%
14	15.718	2260	2266	2277	rVB4	126825	417172	29.29%	4.762%

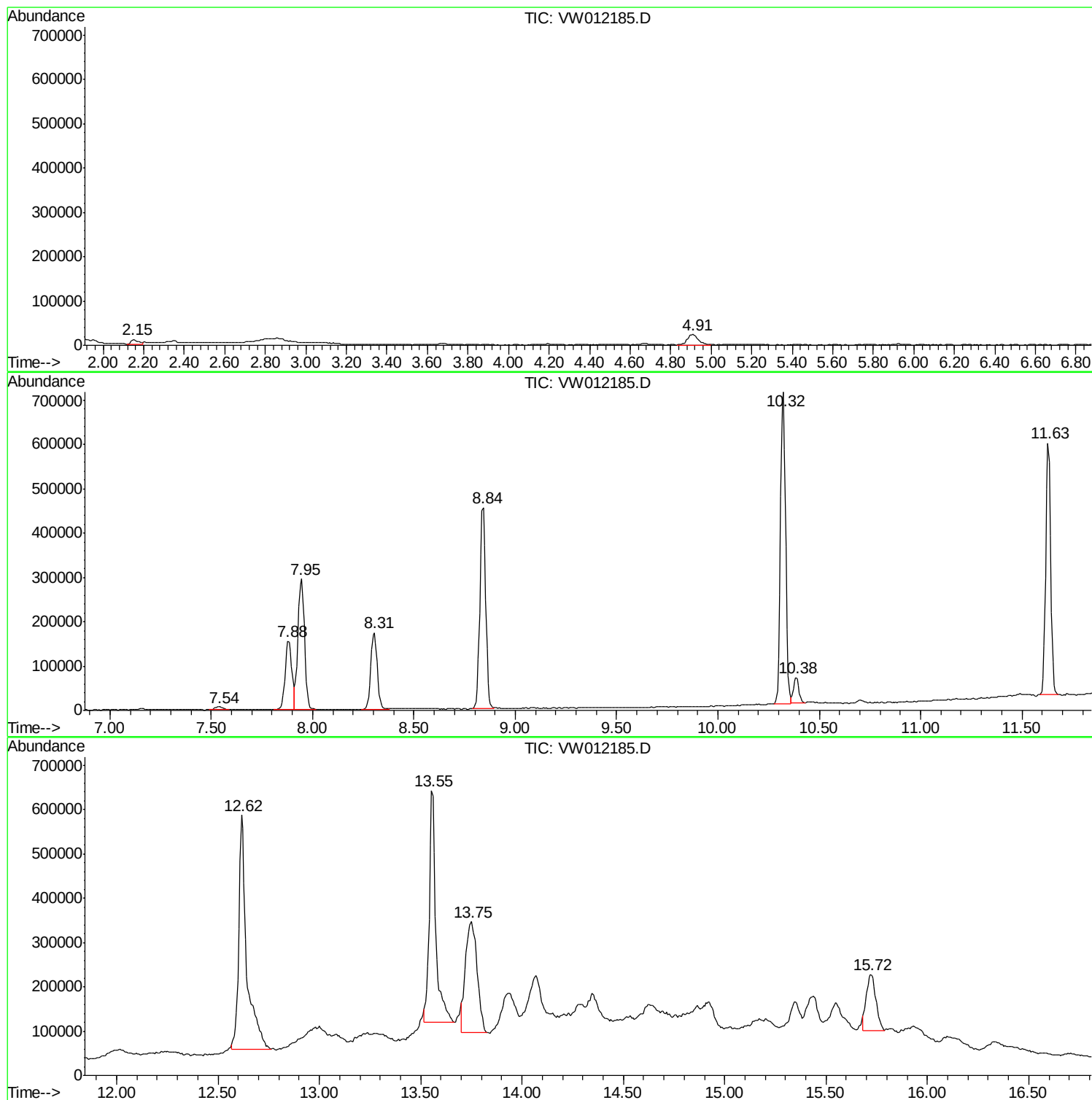
Sum of corrected areas: 8761311

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



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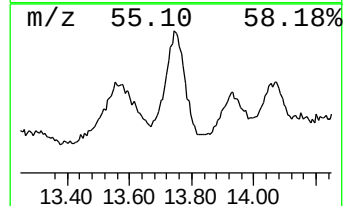
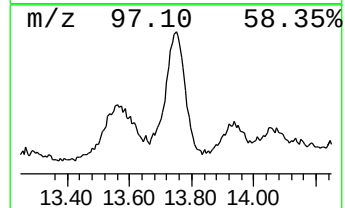
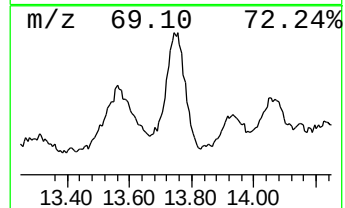
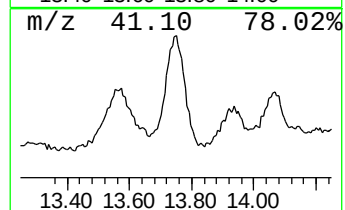
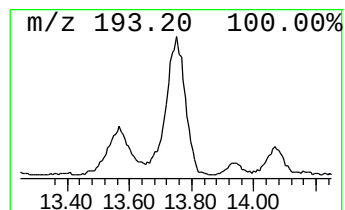
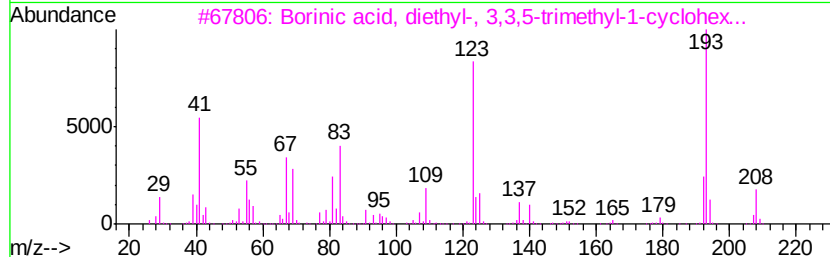
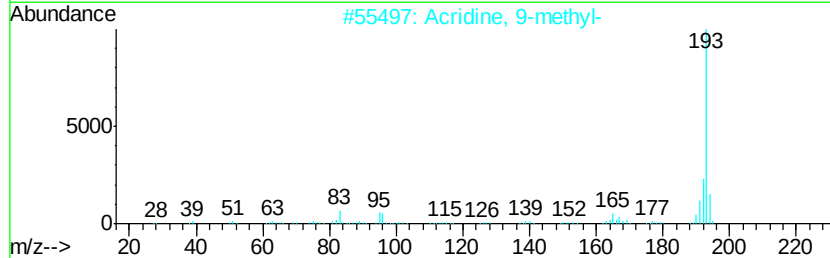
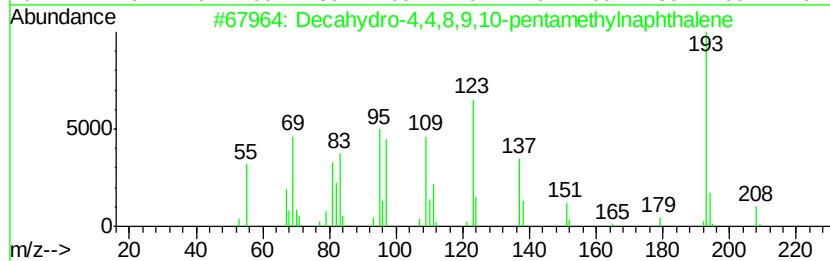
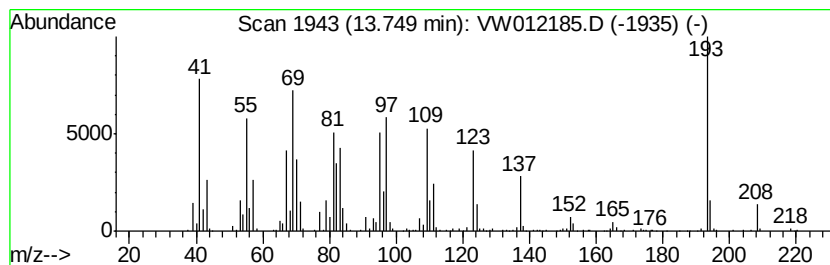
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Peak Number 1 Decahydro-4,4,8,9,10-pentam... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.75	42.35 ug/l	986304	1,4-Dichlorobenzene-d4	13.55
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Decahydro-4,4,8,9,10-pentamethyl...	208 C15H28	080655-44-3 91
2		Acridine, 9-methyl-	193 C14H11N	000611-64-3 30
3		Borinic acid, diethyl-, 3,3,5-tr...	208 C13H25B0	057387-76-5 22
4		2,6-Heptadienal, 2,4-dimethyl-	138 C9H14O	085136-08-9 18
5		[1,1'-Biphenyl]-4-acetonitrile	193 C14H11N	031603-77-7 18



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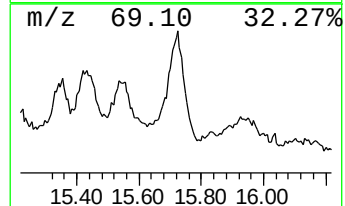
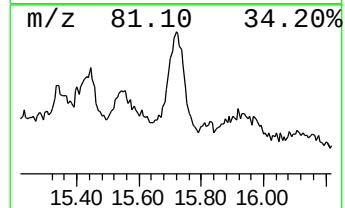
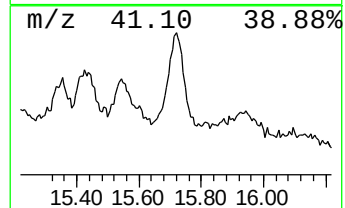
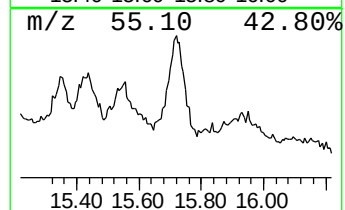
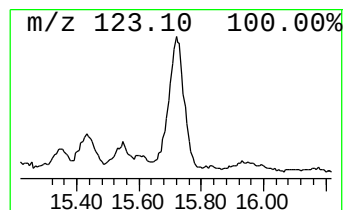
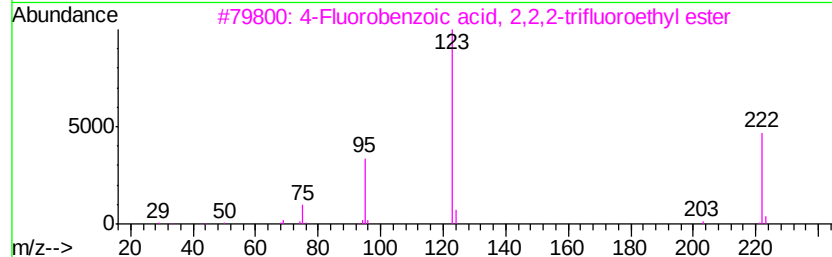
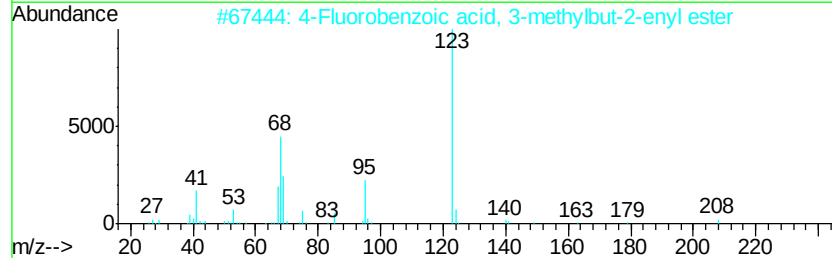
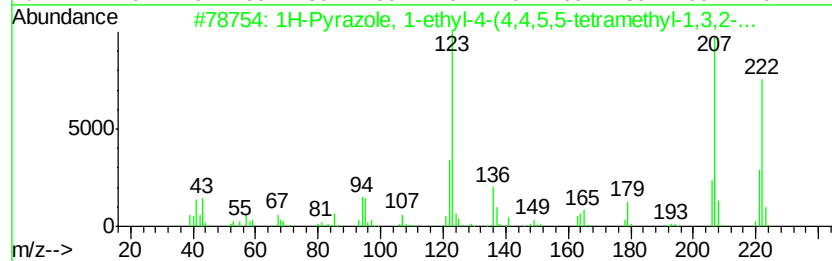
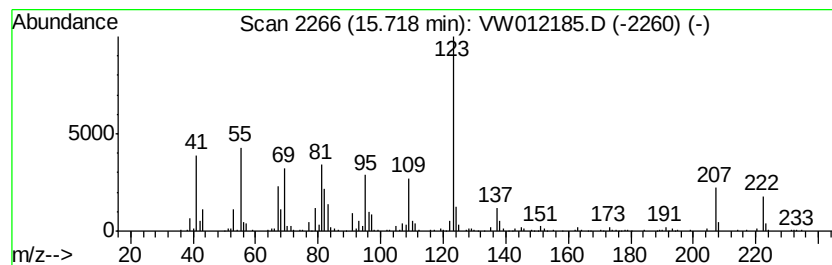
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Peak Number 2 1H-Pyrazole, 1-ethyl-4-(4,4... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.		
15.72	17.91 ug/l	417172	1,4-Dichlorobenzene-d4	13.55		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		1H-Pyrazole, 1-ethyl-4-(4,4,5,5-...	222	C11H19BN2O2	1000319-69-6	53
2		4-Fluorobenzoic acid, 3-methylbu...	208	C12H13F02	1000299-15-1	49
3		4-Fluorobenzoic acid, 2,2,2-trif...	222	C9H6F4O2	1000325-53-2	47
4		1-Trimethylsilylpent-1-en-4-yne	138	C8H14Si	079516-25-9	46
5		4-Fluorobenzoic acid, hex-4-yn-3...	220	C13H13F02	1000299-15-3	46



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TIC Top Hit name	RT	EstConc	Units	Response	---Internal Standard---			
					#	RT	Resp	Conc
Decahydro-4,4,8,9...	13.75	42.4	ug/l	986304	4	13.55	1164390	50.0
1H-Pyrazole, 1-et...	15.72	17.9	ug/l	417172	4	13.55	1164390	50.0