

Data Path : Z:\VOASRV\HPCHEM1\MSVOA W\DATA\VW090819\
 Data File : VW012804.D
 Acq On : 09 Sep 2019 06:51
 Operator : SY/VA
 Sample : K4760-01
 Misc : 5.25G/5ML/MSVOA W/SOIL
 ALS Vial : 45 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 EAST-SIDE-WALL-1

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_W\METHOD\82W090719S.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	2.160	37	42	47	rVB2	6598	10918	10.31%	1.774%
2	2.636	115	120	125	rVB6	3683	6560	6.19%	1.066%
3	2.867	154	158	159	rVB3	1069	1158	1.09%	0.188%
4	2.989	175	178	180	rBV3	1129	1945	1.84%	0.316%
5	3.026	180	184	190	rVB5	3717	8777	8.29%	1.426%
6	3.398	239	245	249	rBV6	1492	3409	3.22%	0.554%
7	3.757	299	304	310	rVB4	741	1371	1.29%	0.223%
8	4.123	352	364	372	rBV3	5433	16749	15.81%	2.722%
9	4.903	484	492	495	rBV6	1837	4609	4.35%	0.749%
10	7.147	856	860	865	rVB5	598	1070	1.01%	0.174%
11	7.891	973	982	986	rBV3	14852	36646	34.60%	5.955%
12	7.946	986	991	1000	rVB	32858	71028	67.06%	11.542%
13	8.305	1042	1050	1060	rBV2	17392	37400	35.31%	6.077%
14	8.842	1130	1138	1147	rBV	41706	85304	80.54%	13.862%
15	9.086	1174	1178	1185	rVB4	3615	7049	6.66%	1.145%
16	10.323	1373	1381	1390	rBV	59406	105910	100.00%	17.210%
17	10.707	1437	1444	1450	rVB3	4734	7827	7.39%	1.272%
18	10.860	1464	1469	1476	rVB2	9288	17583	16.60%	2.857%
19	11.091	1503	1507	1510	rVB4	907	1137	1.07%	0.185%
20	11.628	1589	1595	1603	rBV	45358	77957	73.61%	12.668%
21	12.615	1749	1757	1765	rBV2	29005	53942	50.93%	8.765%
22	13.506	1899	1903	1906	rBV4	929	1383	1.31%	0.225%
23	13.560	1906	1912	1918	rVB	18997	32706	30.88%	5.315%
24	13.652	1922	1927	1931	rVB2	2407	3570	3.37%	0.580%
25	14.066	1990	1995	2002	rVB2	2682	4684	4.42%	0.761%
26	14.524	2063	2070	2076	rVB5	1605	3145	2.97%	0.511%
27	14.725	2099	2103	2106	rVB3	1873	1901	1.79%	0.309%
28	15.152	2171	2173	2177	rBV3	834	1187	1.12%	0.193%
29	15.249	2186	2189	2191	rBV3	967	1249	1.18%	0.203%
30	15.408	2213	2215	2218	rBV3	892	1096	1.03%	0.178%
31	15.572	2241	2242	2248	rVB5	1074	1237	1.17%	0.201%
32	16.182	2338	2342	2345	rBV3	1524	2152	2.03%	0.350%
33	16.401	2375	2378	2381	rVB4	820	1270	1.20%	0.206%
34	16.432	2381	2383	2386	rBV4	1098	1464	1.38%	0.238%

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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

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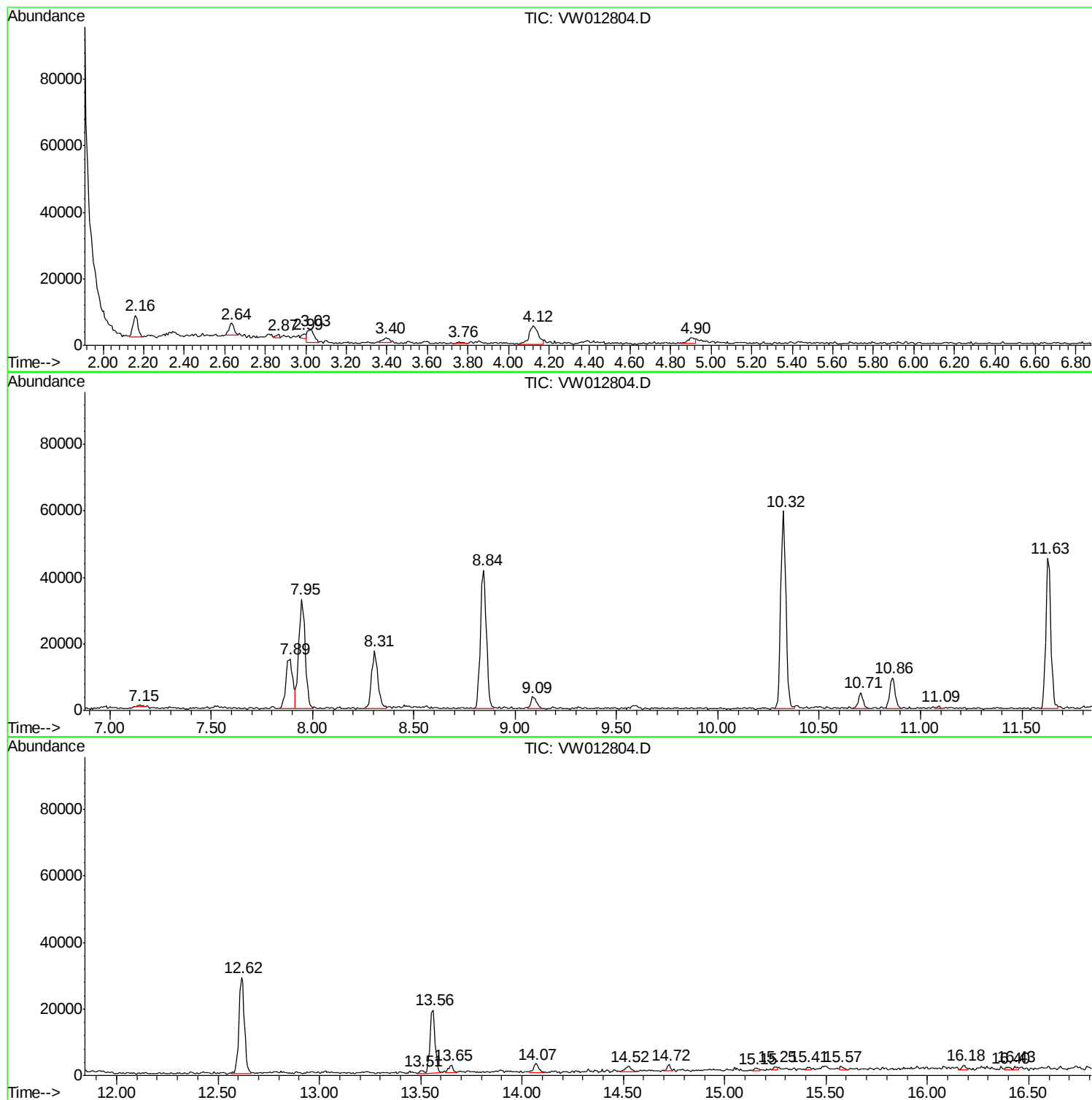
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W090719S.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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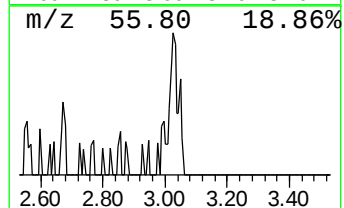
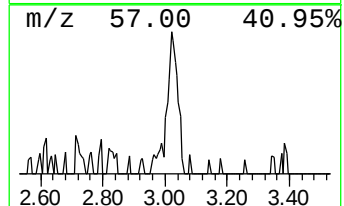
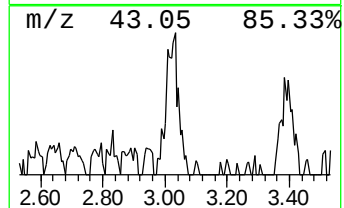
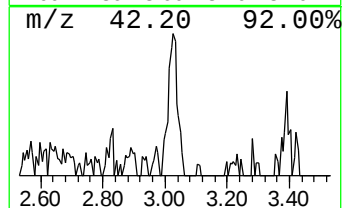
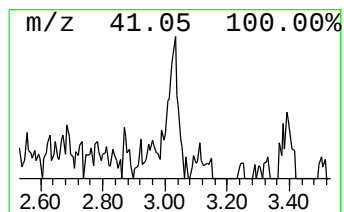
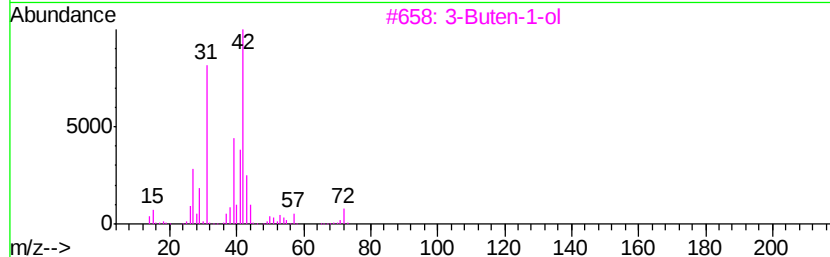
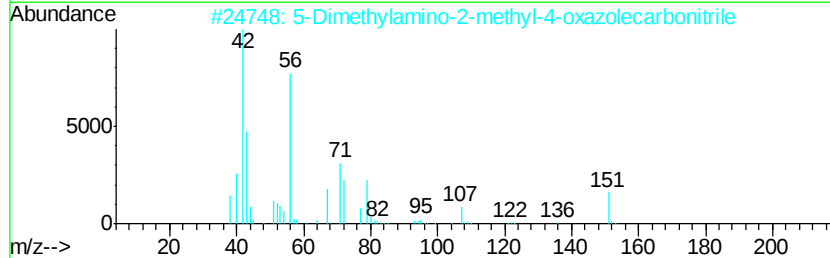
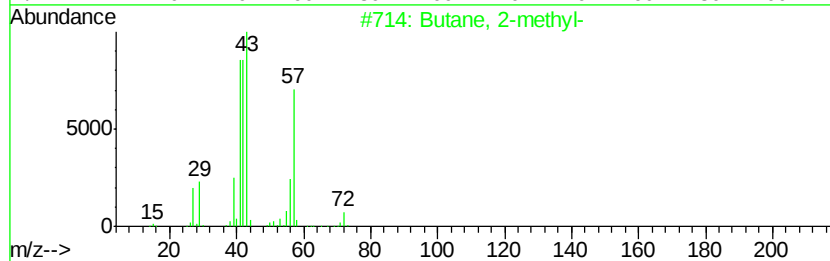
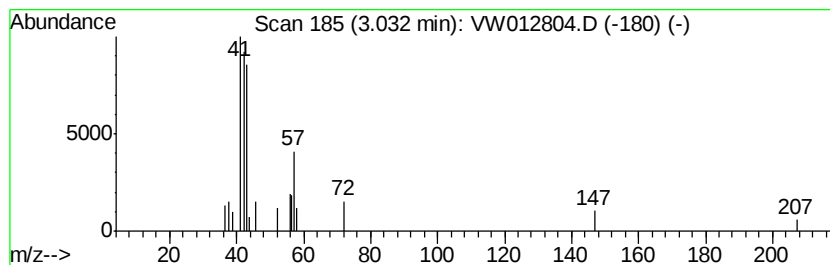
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TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown3.03 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.03	6.18 ug/l	8777	Pentafluorobenzene	7.95

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methyl-	72	C5H12	000078-78-4	35
2		5-Dimethylamino-2-methyl-4-oxazo...	151	C7H9N3O	049837-48-1	17
3		3-Buten-1-ol	72	C4H8O	000627-27-0	9
4		1-Butene	56	C4H8	000106-98-9	9
5		1,3-Dimethyldiaziridine	72	C3H8N2	026177-36-6	9



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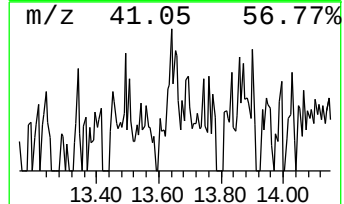
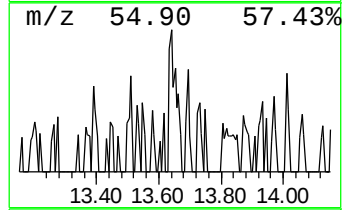
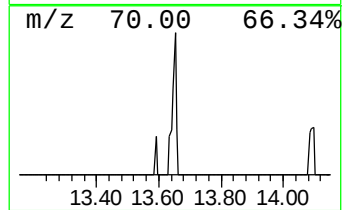
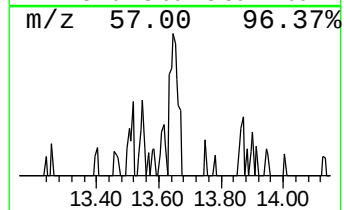
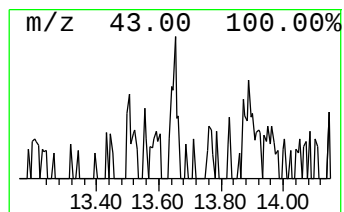
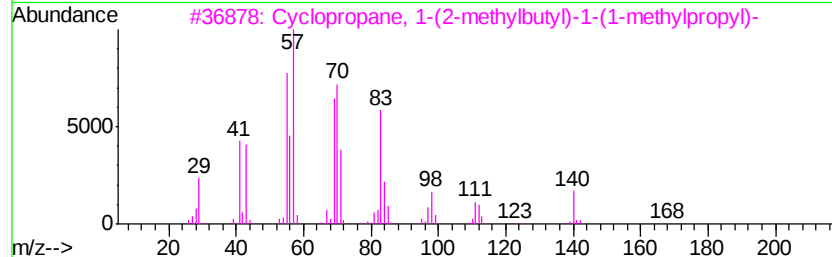
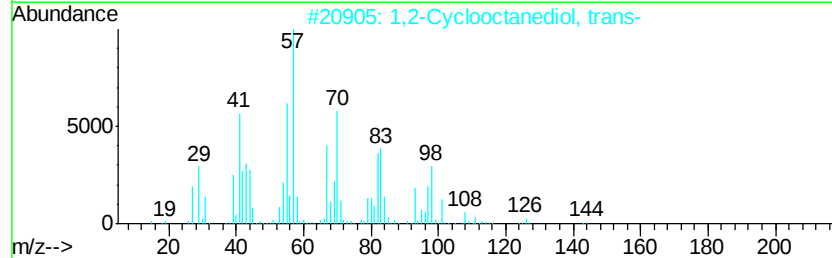
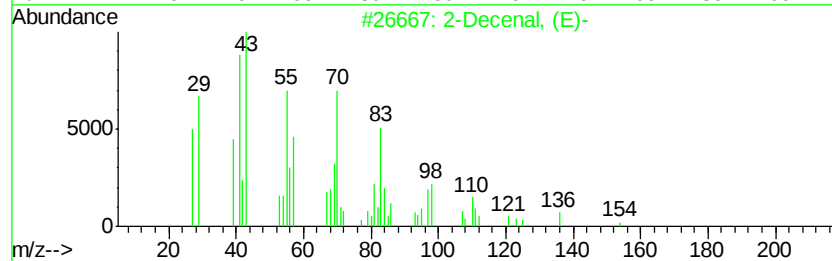
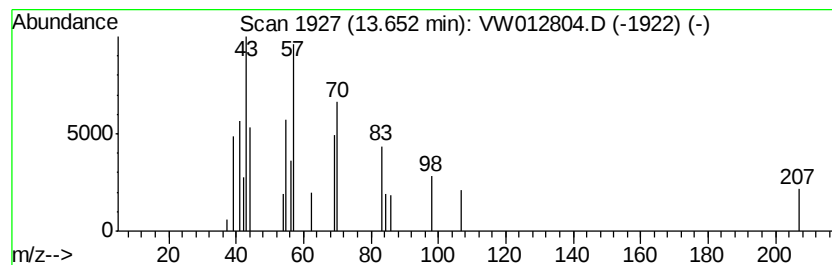
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Peak Number 4 unknown13.65 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.65	5.46 ug/l	3570	1,4-Dichlorobenzene-d4	13.56

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Decenal, (E)-			154	C10H18O	003913-81-3	27
2	1,2-Cyclooctanediol, trans-			144	C8H16O2	042565-22-0	25
3	Cyclopropane, 1-(2-methylbutyl)-...			168	C12H24	064723-36-0	16
4	2-Octanol, trifluoroacetate			226	C10H17F3O2	1000352-37-1	14
5	1-Pentene, 2,3-dimethyl-			98	C7H14	003404-72-6	14



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown3.03	3.03	6.2	ug/l	8777	1	7.95	71028	50.0
unknown13.65	13.65	5.5	ug/l	3570	4	13.56	32706	50.0