

Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD121121\
 Data File : VD071279.D
 Acq On : 11 Dec 2021 19:37
 Operator : VA/SY
 Sample : M5017-09
 Misc : 5.01G/5.00ml/MSVOA_D/SOIL
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 TP1

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\82D112921S.M
 Title : SW846 8260

Signal : TIC: VD071279.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.232	49	53	55	rVV	4496	5293	1.54%	0.270%
2	2.644	115	123	127	rVB3	8087	20288	5.89%	1.036%
3	3.138	205	207	215	rVB3	8877	16846	4.89%	0.860%
4	4.785	485	487	491	rVB	3299	4614	1.34%	0.236%
5	4.838	491	496	498	rBV	3885	6663	1.93%	0.340%
6	4.961	510	517	518	rBV3	7389	11811	3.43%	0.603%
7	5.191	552	556	560	rVB4	2715	4736	1.37%	0.242%
8	5.979	684	690	692	rBV2	2131	3468	1.01%	0.177%
9	6.256	734	737	741	rVB2	2363	4047	1.17%	0.207%
10	7.197	889	897	904	rVB3	10279	29054	8.43%	1.483%
11	7.691	976	981	983	rBV	3357	4456	1.29%	0.227%
12	7.914	1010	1019	1024	rBV2	39535	98387	28.56%	5.022%
13	7.973	1024	1029	1039	rVV	68594	150239	43.61%	7.669%
14	8.173	1060	1063	1066	rBV2	1966	3505	1.02%	0.179%
15	8.332	1081	1090	1096	rBV2	47930	109679	31.83%	5.599%
16	8.861	1173	1180	1187	rBV	101432	201925	58.61%	10.308%
17	9.149	1225	1229	1232	rBV	2268	3502	1.02%	0.179%
18	10.332	1423	1430	1438	rBV	149674	281422	81.68%	14.366%
19	11.443	1616	1619	1623	rVB	3575	3934	1.14%	0.201%
20	11.638	1643	1652	1662	rBV	176561	309879	89.94%	15.819%
21	12.079	1725	1727	1731	rBV2	2810	3739	1.09%	0.191%
22	12.620	1812	1819	1827	rBV	148230	258021	74.89%	13.171%
23	12.873	1859	1862	1867	rBV2	3633	4713	1.37%	0.241%
24	12.943	1870	1874	1879	rVB6	5334	8954	2.60%	0.457%
25	13.173	1909	1913	1916	rBV4	3768	5231	1.52%	0.267%
26	13.255	1923	1927	1932	rBV5	4379	8149	2.37%	0.416%
27	13.567	1972	1980	1988	rVB	195818	344528	100.00%	17.587%
28	13.879	2028	2033	2037	rBV3	5325	8230	2.39%	0.420%
29	13.932	2039	2042	2044	rBV2	2653	3858	1.12%	0.197%
30	14.908	2205	2208	2212	rVB2	2403	3587	1.04%	0.183%
31	15.084	2235	2238	2242	rVV3	2000	3468	1.01%	0.177%
32	15.384	2281	2289	2294	rVB5	11134	24473	7.10%	1.249%
33	15.931	2380	2382	2385	rBV3	3788	4032	1.17%	0.206%
34	16.137	2414	2417	2420	rBV2	3494	4214	1.22%	0.215%

Sum of corrected areas: 1958945

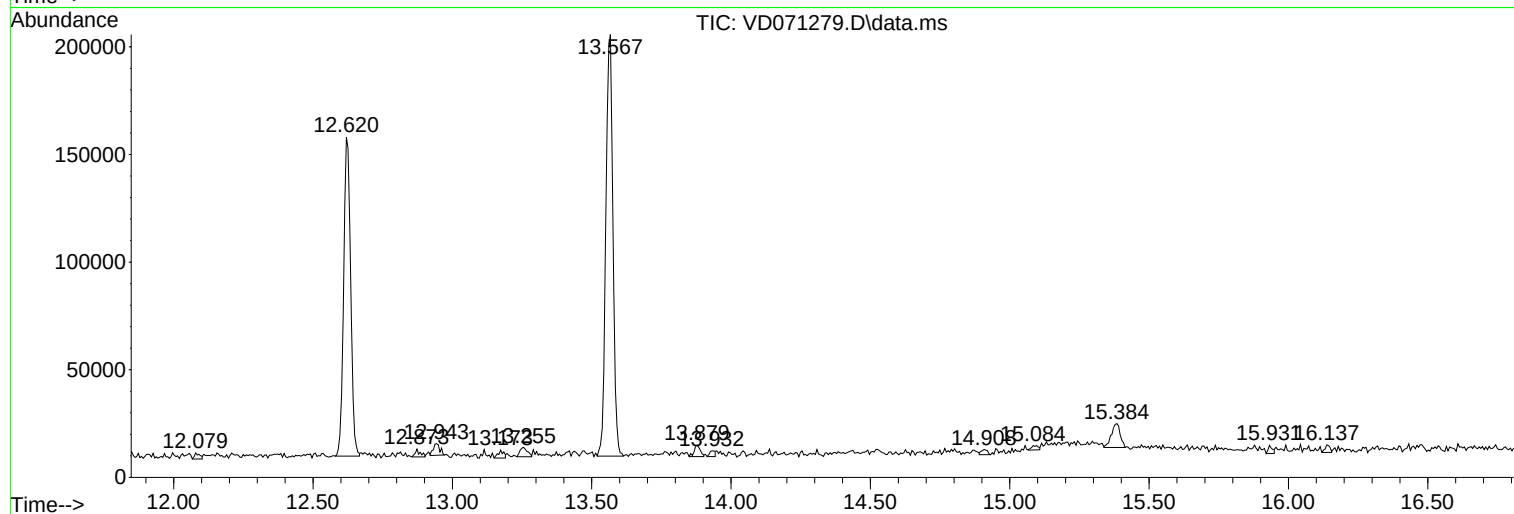
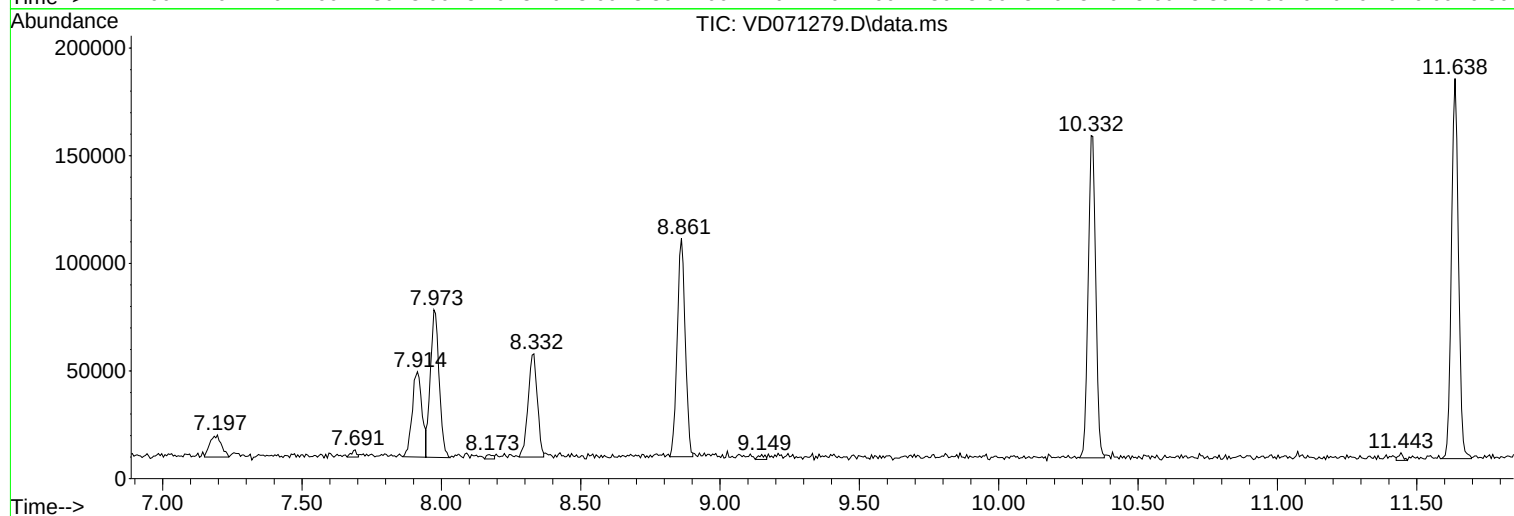
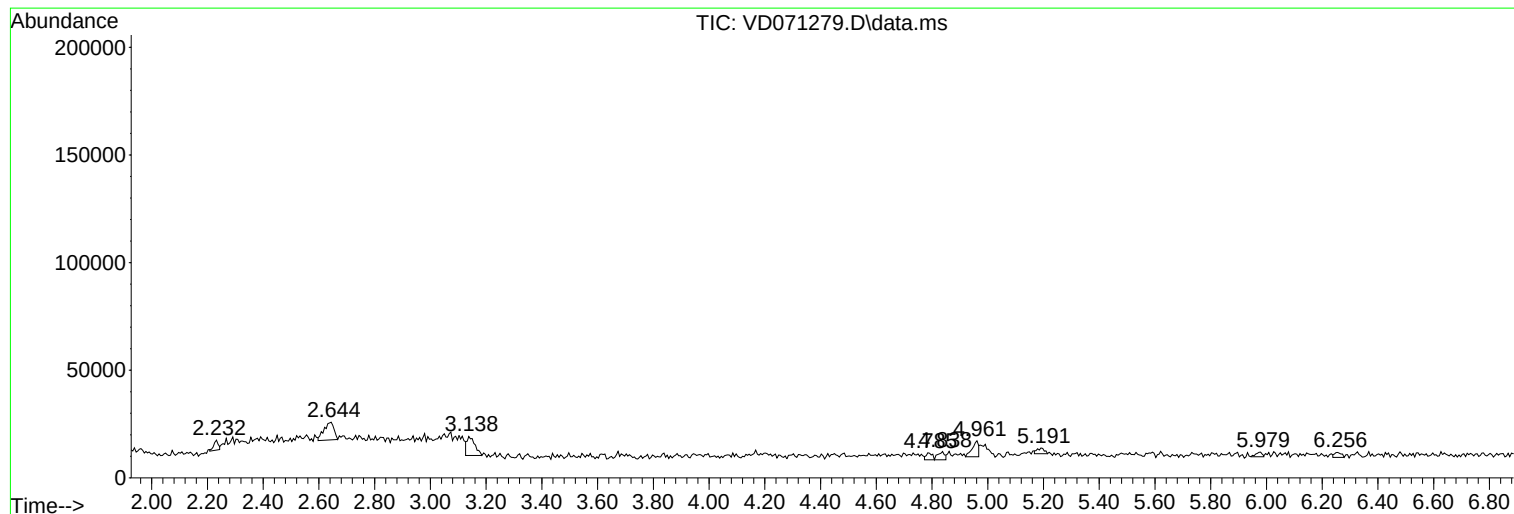
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D112921S.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P



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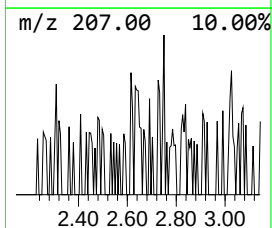
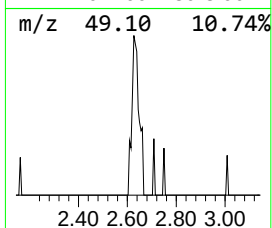
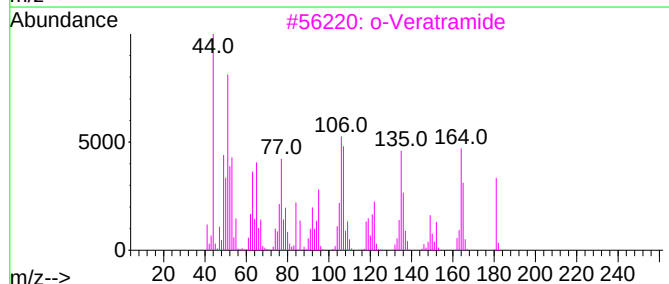
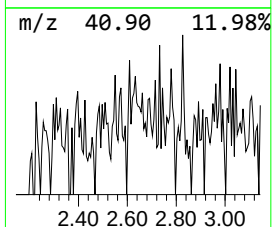
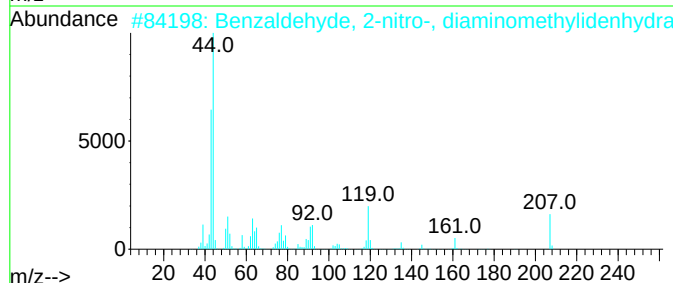
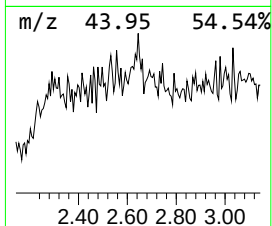
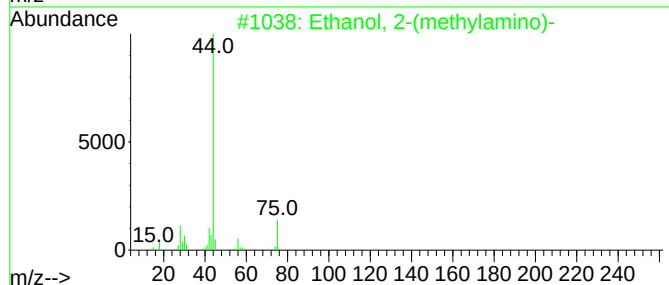
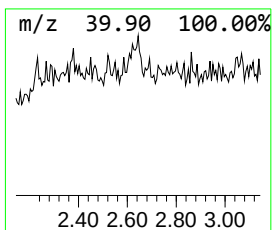
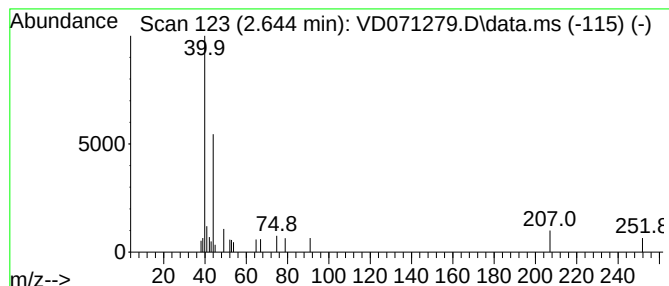
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D112921S.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 unknown2.644 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.644	6.75 ug/l	20288	Pentafluorobenzene	7.973

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-(methylamino)-	75	C3H9NO	000109-83-1	25
2			Benzaldehyde, 2-nitro-, diaminom...	207	C8H9N5O2	102632-31-5	9
3			o-Veratramide	181	C9H11NO3	001521-39-7	9
4			5-Aminoisoxazole	84	C3H4N2O	014678-05-8	9
5			Acetonitrile-d3	44	C2D3N	002206-26-0	4



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown2.644	2.644	6.8	ug/l	20288	1	7.973	150239	50.0