

Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW082622\
 Data File : VW024371.D
 Acq On : 27 Aug 2022 00:01
 Operator : SY/VA
 Sample : N4344-11
 Misc : 5.60g/10mL/MSVOA_W/SOIL
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 DBVR0

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0 % 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\SFAMWLM082622SMA.M
 Title : SFAM01.0

Signal : TIC: VW024371.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.251	47	57	58	rBV9	12997	33847	4.67%	0.558%
2	2.343	66	72	79	rBV	129257	226847	31.31%	3.741%
3	2.873	153	159	168	rVB	77812	161393	22.28%	2.662%
4	3.458	252	255	257	rBV4	1680	1933	0.27%	0.032%
5	3.641	284	285	287	rBV2	1430	1389	0.19%	0.023%
6	3.721	295	298	300	rBV4	1595	1488	0.21%	0.025%
7	3.903	326	328	330	rBV3	1899	1907	0.26%	0.031%
8	4.007	335	345	356	rVV2	61141	187554	25.89%	3.093%
9	4.105	359	361	364	rVB4	1705	1634	0.23%	0.027%
10	4.281	388	390	394	rVB5	2092	1889	0.26%	0.031%
11	4.355	401	402	405	rVB3	3142	2256	0.31%	0.037%
12	4.397	405	409	411	rBV5	2505	4731	0.65%	0.078%
13	4.483	421	423	428	rVB6	3170	4833	0.67%	0.080%
14	4.623	441	446	447	rBV5	1907	3008	0.42%	0.050%
15	4.915	484	494	502	rBV5	12556	42074	5.81%	0.694%
16	5.013	508	510	513	rVB4	1810	1605	0.22%	0.026%
17	5.043	513	515	516	rBV2	2322	1856	0.26%	0.031%
18	5.098	520	524	525	rBV4	2195	2351	0.32%	0.039%
19	5.281	551	554	555	rBV3	1807	2076	0.29%	0.034%
20	5.379	568	570	577	rVB8	2163	4091	0.56%	0.067%
21	5.562	598	600	604	rVB4	2041	1870	0.26%	0.031%
22	5.860	646	649	652	rBV4	1249	1862	0.26%	0.031%
23	5.927	656	660	665	rVB7	3283	7620	1.05%	0.126%
24	6.183	701	702	708	rBV6	1303	2062	0.28%	0.034%
25	6.409	737	739	742	rVB4	1599	1637	0.23%	0.027%
26	6.452	742	746	749	rBV5	1429	1971	0.27%	0.033%
27	6.507	753	755	757	rVV3	1828	1564	0.22%	0.026%
28	6.574	761	766	770	rVB8	1802	3384	0.47%	0.056%
29	6.629	773	775	778	rVB4	1388	1409	0.19%	0.023%
30	6.793	797	802	803	rBV5	1157	1571	0.22%	0.026%
31	6.866	812	814	816	rBV3	1504	1680	0.23%	0.028%
32	7.080	841	849	854	rVV2	18268	47746	6.59%	0.787%
33	7.147	855	860	871	rVV	29037	82844	11.44%	1.366%
34	7.537	920	924	927	rVV6	1604	2403	0.33%	0.040%
35	7.647	933	942	952	rVB	122241	296943	40.99%	4.897%
36	7.842	971	974	975	rVB3	1486	1487	0.21%	0.025%

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Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM082622SMA.M
 Title : SFAM01.0

37	7.884	978	981	983	rVB4	1743	1494	0.21%	0.025%
38	7.927	983	988	990	rBV6	1565	2817	0.39%	0.046%
39	8.073	1009	1012	1014	rVB4	1738	1999	0.28%	0.033%
40	8.140	1020	1023	1024	rBV3	1314	1576	0.22%	0.026%
41	8.268	1032	1044	1058	rVV2	232906	724404	100.00%	11.948%
42	8.567	1091	1093	1097	rVB5	3105	2644	0.36%	0.044%
43	8.677	1107	1111	1115	rBV7	2438	3220	0.44%	0.053%
44	8.762	1123	1125	1128	rVB4	1575	1315	0.18%	0.022%
45	8.842	1128	1138	1147	rBV	280900	542745	74.92%	8.951%
46	9.134	1185	1186	1190	rBV4	1589	1697	0.23%	0.028%
47	9.165	1190	1191	1195	rBV4	1321	1505	0.21%	0.025%
48	9.274	1200	1209	1217	rBV	159297	330854	45.67%	5.457%
49	9.396	1227	1229	1233	rBV5	1635	1614	0.22%	0.027%
50	9.530	1249	1251	1252	rBV2	1643	1486	0.21%	0.025%
51	9.756	1286	1288	1290	rVB3	1904	1610	0.22%	0.027%
52	9.774	1290	1291	1294	rBV3	1961	1972	0.27%	0.033%
53	9.890	1307	1310	1313	rVB5	2132	2047	0.28%	0.034%
54	9.969	1322	1323	1325	rBV2	1851	1549	0.21%	0.026%
55	10.030	1325	1333	1342	rVB	72406	138938	19.18%	2.291%
56	10.195	1357	1360	1363	rBV5	1863	2868	0.40%	0.047%
57	10.323	1371	1381	1391	rBV	325221	573806	79.21%	9.464%
58	10.579	1416	1423	1430	rVB	52460	92185	12.73%	1.520%
59	10.670	1435	1438	1439	rBV3	1345	1618	0.22%	0.027%
60	10.701	1439	1443	1446	rVV4	4520	6624	0.91%	0.109%
61	10.859	1461	1469	1475	rVV2	151544	267077	36.87%	4.405%
62	10.927	1475	1480	1489	rVB	72289	127523	17.60%	2.103%
63	11.073	1498	1504	1509	rVB9	7148	14621	2.02%	0.241%
64	11.396	1554	1557	1559	rBV4	2256	2265	0.31%	0.037%
65	11.433	1560	1563	1569	rVB8	3383	6379	0.88%	0.105%
66	11.628	1587	1595	1602	rBV	418894	702379	96.96%	11.584%
67	11.841	1627	1630	1633	rVB5	2306	2345	0.32%	0.039%
68	12.176	1678	1685	1687	rBV8	2468	4339	0.60%	0.072%
69	12.341	1704	1712	1714	rBV8	1758	3981	0.55%	0.066%
70	12.463	1729	1732	1733	rBV3	1522	1395	0.19%	0.023%
71	12.603	1750	1755	1760	rVV9	5437	10164	1.40%	0.168%
72	12.688	1762	1769	1777	rBV	129400	217959	30.09%	3.595%
73	12.847	1793	1795	1798	rBV3	1331	1306	0.18%	0.022%
74	12.944	1809	1811	1813	rVB3	2698	1987	0.27%	0.033%
75	13.030	1823	1825	1828	rBV4	1670	1957	0.27%	0.032%
76	13.402	1882	1886	1890	rBV7	2968	5239	0.72%	0.086%

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Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM082622SMA.M
 Title : SFAM01.0

77	13.554	1905	1911	1919	rVV	384778	594670	82.09%	9.808%
78	13.646	1924	1926	1931	rVB5	4327	5763	0.80%	0.095%
79	13.694	1931	1934	1935	rBV3	1937	1942	0.27%	0.032%
80	13.743	1940	1942	1944	rBV3	2020	1803	0.25%	0.030%
81	13.847	1950	1959	1972	rVB	258915	442733	61.12%	7.302%
82	14.060	1991	1994	1999	rVV7	3093	4649	0.64%	0.077%
83	14.219	2016	2020	2022	rBV5	1321	1738	0.24%	0.029%
84	14.310	2033	2035	2036	rBV2	3806	2936	0.41%	0.048%
85	14.536	2068	2072	2077	rVB8	2958	6084	0.84%	0.100%
86	14.578	2077	2079	2080	rBV2	2470	2467	0.34%	0.041%
87	14.694	2096	2098	2100	rBV3	1812	1842	0.25%	0.030%
88	14.719	2100	2102	2105	rVV4	2366	3449	0.48%	0.057%
89	14.993	2143	2147	2148	rBV4	1238	1367	0.19%	0.023%
90	15.255	2189	2190	2192	rVB2	2477	1740	0.24%	0.029%
91	15.279	2192	2194	2196	rBV2	1978	2239	0.31%	0.037%
92	15.426	2214	2218	2224	rBV6	5632	9652	1.33%	0.159%
93	15.651	2253	2255	2258	rVB4	1238	1331	0.18%	0.022%
94	15.773	2273	2275	2283	rVB8	1773	2786	0.38%	0.046%
95	15.834	2283	2285	2288	rBV4	1359	1654	0.23%	0.027%
96	15.920	2294	2299	2300	rBV5	1195	1809	0.25%	0.030%
97	16.054	2319	2321	2326	rVV6	1750	1910	0.26%	0.032%
98	16.255	2351	2354	2357	rBV5	1591	1941	0.27%	0.032%
99	16.279	2357	2358	2363	rVV5	2137	2358	0.33%	0.039%
100	16.444	2380	2385	2391	rBV10	1896	4037	0.56%	0.067%

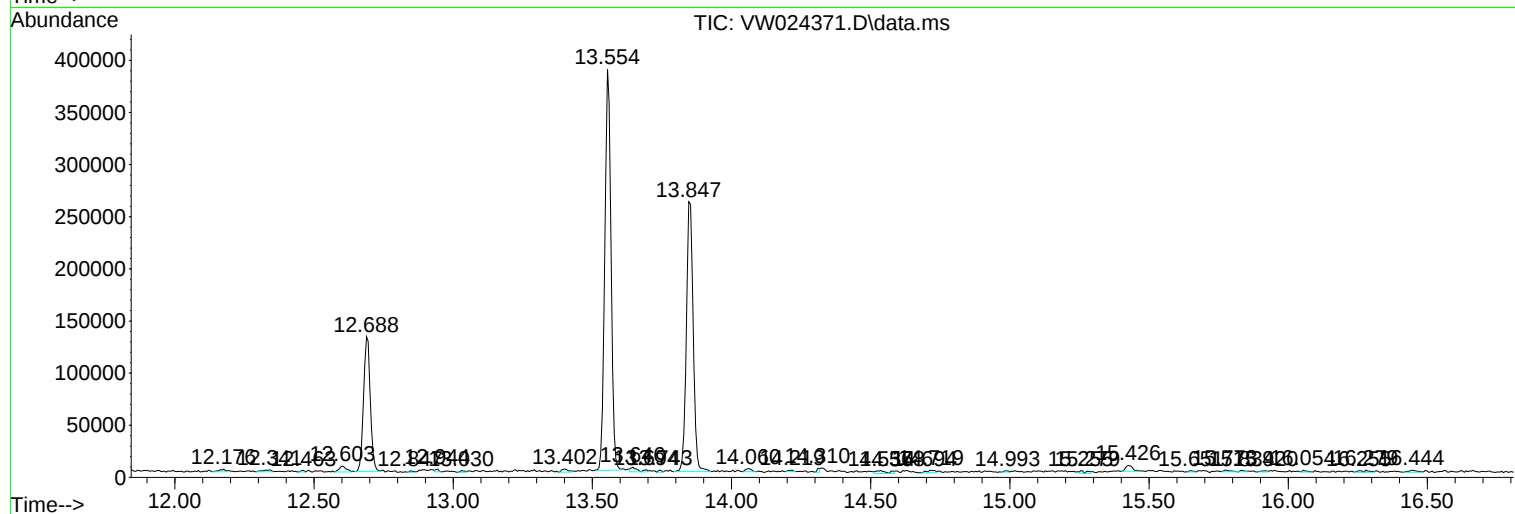
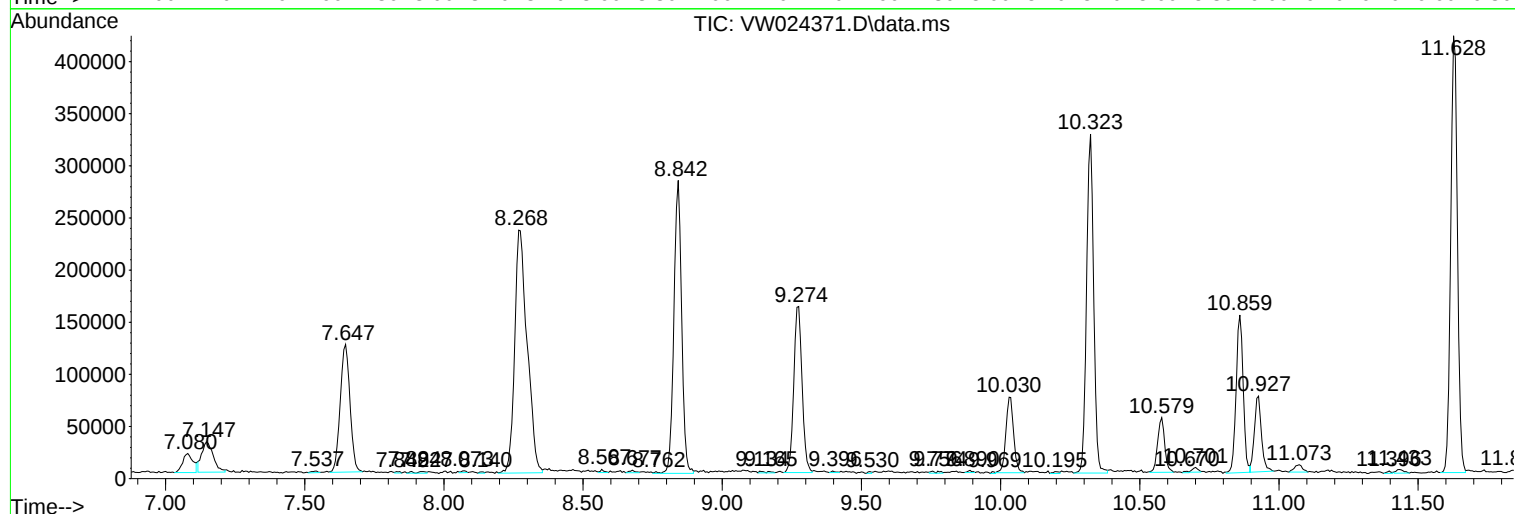
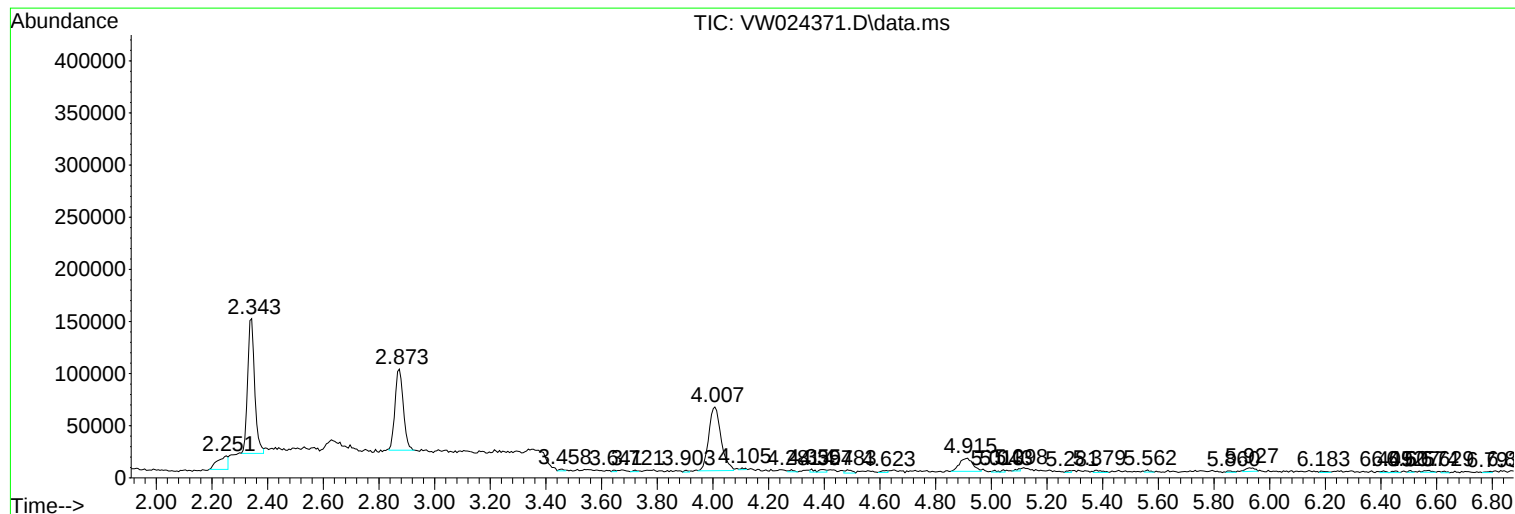
Sum of corrected areas: 6063218

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Instrument :
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DBVR0

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM082622SMA.M
Quant Title : SFAM01.0

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



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Instrument :
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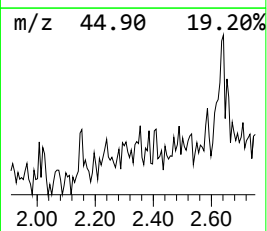
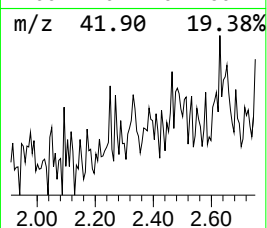
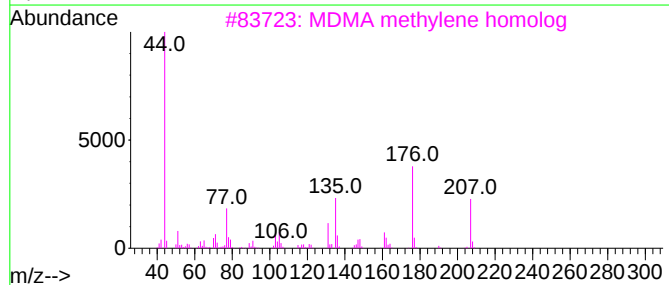
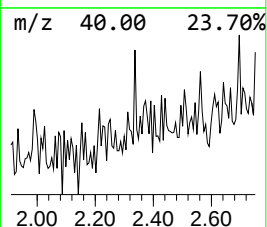
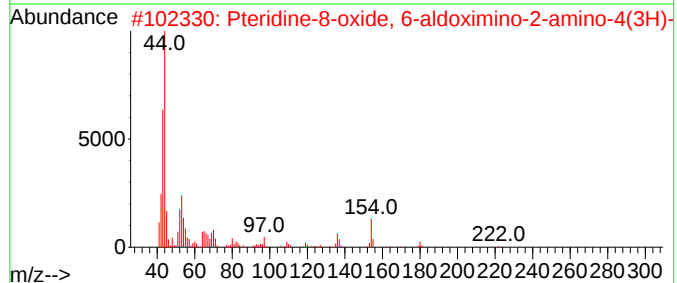
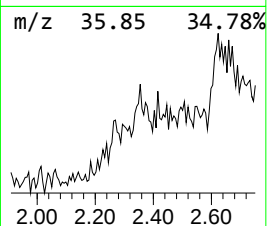
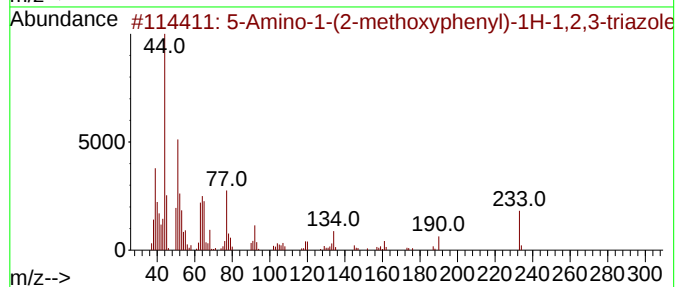
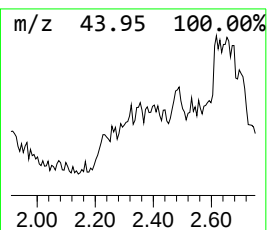
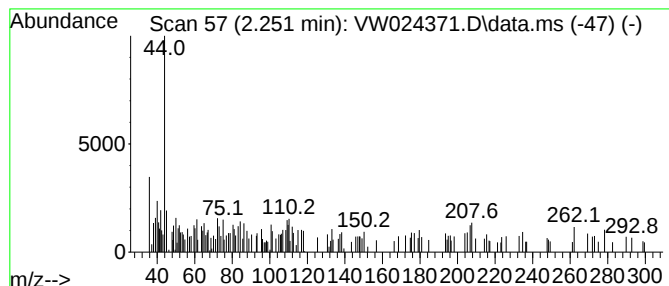
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM082622SMA.M
Quant Title : SFAM01.0

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

Peak Number 1 5-Amino-1-(2-methoxyphenyl)... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.251	1.56 ug/L	33847	1,4-Difluorobenzene	8.842

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			5-Amino-1-(2-methoxyphenyl)-1H-1...	233	C10H11N5O2	1000387-16-8	53
2			Pteridine-8-oxide, 6-aldoximino-...	222	C7H6N6O3	023663-23-2	38
3			MDMA methylene homolog	207	C12H17NO2	1010386-32-2	25
4			Benzaldehyde, 2-nitro-, diaminom...	207	C8H9N5O2	102632-31-5	25
5			Acetamide, 2,2-dichloro-	127	C2H3Cl2NO	000683-72-7	22



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
5-Amino-1-(2-me...	2.251	1.6	ug/L	33847	1	8.842	542745	25.0