

Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW042822\
 Data File : VW022677.D
 Acq On : 28 Apr 2022 21:58
 Operator : SY/VA
 Sample : N2603-13
 Misc : 8.86g/10mL/MSVOA_W/SOIL
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 EW494

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\SFAMWLM040722SMA.M
 Title : SFAM01.0

Signal : TIC: VW022677.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.355	69	74	80	rVB	51680	84204	21.10%	3.695%
2	3.276	223	225	227	rBV4	942	834	0.21%	0.037%
3	3.477	256	258	260	rBV3	1061	710	0.18%	0.031%
4	3.593	275	277	278	rBV2	1341	727	0.18%	0.032%
5	3.672	288	290	292	rBV3	839	684	0.17%	0.030%
6	3.800	310	311	314	rVB3	900	663	0.17%	0.029%
7	4.013	338	346	358	rVB2	38011	107910	27.04%	4.735%
8	4.251	383	385	389	rBV5	1218	1781	0.45%	0.078%
9	4.611	442	444	447	rVB4	1108	973	0.24%	0.043%
10	4.678	451	455	457	rBV4	1297	1559	0.39%	0.068%
11	4.794	472	474	478	rBV4	1031	1284	0.32%	0.056%
12	5.092	521	523	528	rBV5	1598	2215	0.56%	0.097%
13	5.129	528	529	533	rVB4	1276	1567	0.39%	0.069%
14	5.312	556	559	561	rVB4	820	695	0.17%	0.030%
15	5.367	565	568	569	rBV2	907	852	0.21%	0.037%
16	5.531	591	595	597	rVB3	1366	1547	0.39%	0.068%
17	5.556	597	599	603	rBV5	910	1425	0.36%	0.063%
18	5.598	603	606	609	rBV3	1415	1876	0.47%	0.082%
19	5.726	625	627	630	rVB3	1226	944	0.24%	0.041%
20	5.757	630	632	633	rBV2	1116	875	0.22%	0.038%
21	5.806	637	640	643	rVB4	1080	1403	0.35%	0.062%
22	5.928	653	660	661	rBV6	2388	4400	1.10%	0.193%
23	6.025	674	676	680	rBV5	785	1016	0.25%	0.045%
24	6.117	687	691	693	rBV2	1153	1630	0.41%	0.072%
25	6.196	702	704	708	rVB5	1032	823	0.21%	0.036%
26	6.312	720	723	726	rVB3	1036	1073	0.27%	0.047%
27	6.458	744	747	748	rBV3	1069	1109	0.28%	0.049%
28	6.531	757	759	760	rBV2	835	769	0.19%	0.034%
29	6.769	796	798	800	rBV3	732	810	0.20%	0.036%
30	6.982	831	833	836	rBV2	1204	1312	0.33%	0.058%
31	7.080	840	849	866	rBV	31649	100422	25.17%	4.406%
32	7.385	896	899	901	rBV4	861	776	0.19%	0.034%
33	7.653	934	943	952	rBV2	81159	205681	51.54%	9.025%
34	7.750	956	959	960	rVB3	896	793	0.20%	0.035%
35	7.836	972	973	975	rBV2	1245	916	0.23%	0.040%
36	7.891	980	982	983	rBV	939	710	0.18%	0.031%

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Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM040722SMA.M
 Title : SFAM01.0

37	8.281	1035	1046	1058	rBV2	116952	399051	100.00%	17.510%
38	8.842	1130	1138	1147	rVV	117808	235117	58.92%	10.317%
39	8.988	1160	1162	1165	rVB4	864	844	0.21%	0.037%
40	9.073	1174	1176	1177	rBV	1730	1190	0.30%	0.052%
41	9.165	1189	1191	1195	rBV5	1104	1216	0.30%	0.053%
42	9.274	1200	1209	1218	rBV	106135	217427	54.49%	9.540%
43	9.463	1239	1240	1242	rBV2	1121	892	0.22%	0.039%
44	9.531	1249	1251	1255	rBB5	981	967	0.24%	0.042%
45	9.659	1270	1272	1273	rBV2	928	681	0.17%	0.030%
46	9.793	1293	1294	1297	rVB2	1269	1208	0.30%	0.053%
47	9.829	1297	1300	1303	rBV4	1275	1595	0.40%	0.070%
48	9.866	1303	1306	1307	rVV3	1383	1401	0.35%	0.061%
49	9.884	1307	1309	1314	rVV6	2736	3905	0.98%	0.171%
50	9.976	1323	1324	1325	rVV	1288	658	0.16%	0.029%
51	10.037	1327	1334	1343	rVB2	28489	53626	13.44%	2.353%
52	10.323	1374	1381	1388	rBV	84282	147795	37.04%	6.485%
53	10.500	1405	1410	1411	rBV4	1230	1785	0.45%	0.078%
54	10.579	1417	1423	1431	rBV	14621	27049	6.78%	1.187%
55	10.701	1439	1443	1448	rBV3	5304	7120	1.78%	0.312%
56	10.927	1474	1480	1490	rBV	85419	156780	39.29%	6.879%
57	11.219	1526	1528	1530	rVB3	1215	964	0.24%	0.042%
58	11.335	1545	1547	1550	rBV4	1365	1749	0.44%	0.077%
59	11.366	1550	1552	1555	rVV3	957	1183	0.30%	0.052%
60	11.433	1561	1563	1566	rVB4	1085	897	0.22%	0.039%
61	11.500	1572	1574	1576	rVV3	1026	770	0.19%	0.034%
62	11.518	1576	1577	1581	rVB3	1201	1156	0.29%	0.051%
63	11.628	1588	1595	1604	rBV	72340	134332	33.66%	5.894%
64	11.835	1627	1629	1631	rVV3	1325	1031	0.26%	0.045%
65	11.920	1639	1643	1646	rBV7	897	1500	0.38%	0.066%
66	12.054	1663	1665	1668	rVB2	865	798	0.20%	0.035%
67	12.128	1676	1677	1679	rVV2	1172	825	0.21%	0.036%
68	12.164	1682	1683	1685	rVV2	1552	1307	0.33%	0.057%
69	12.341	1710	1712	1719	rVB6	1065	1491	0.37%	0.065%
70	12.445	1725	1729	1730	rBV3	958	1193	0.30%	0.052%
71	12.603	1749	1755	1762	rVB4	6529	12943	3.24%	0.568%
72	12.689	1762	1769	1776	rBV	113071	189985	47.61%	8.336%
73	12.914	1801	1806	1807	rBV4	1315	1688	0.42%	0.074%
74	13.085	1831	1834	1838	rBV5	1027	1486	0.37%	0.065%
75	13.262	1861	1863	1865	rBV2	1074	691	0.17%	0.030%
76	13.554	1906	1911	1918	rBV	25395	43280	10.85%	1.899%

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 Title : SFAM01.0

77	13.713	1935	1937	1939	rVB3	1403	916	0.23%	0.040%
78	13.737	1939	1941	1944	rBV4	618	879	0.22%	0.039%
79	13.853	1949	1960	1968	rBV	26285	52869	13.25%	2.320%
80	13.987	1981	1982	1985	rBV2	884	882	0.22%	0.039%
81	14.018	1985	1987	1990	rVV4	1075	898	0.23%	0.039%
82	14.060	1991	1994	1999	rVV3	3172	4900	1.23%	0.215%
83	14.133	2002	2006	2010	rVB5	855	1214	0.30%	0.053%
84	14.225	2020	2021	2025	rBV3	1246	907	0.23%	0.040%
85	14.322	2035	2037	2043	rVB6	1862	2155	0.54%	0.095%
86	14.414	2051	2052	2055	rVB3	1124	744	0.19%	0.033%
87	14.609	2081	2084	2089	rBV5	1027	2156	0.54%	0.095%
88	14.847	2121	2123	2125	rVB3	893	710	0.18%	0.031%
89	14.920	2132	2135	2137	rBV3	917	943	0.24%	0.041%
90	15.091	2162	2163	2165	rBV2	1437	1184	0.30%	0.052%
91	15.176	2175	2177	2180	rBV3	946	867	0.22%	0.038%
92	15.462	2221	2224	2225	rBV3	928	1005	0.25%	0.044%
93	15.615	2247	2249	2250	rBV2	1042	731	0.18%	0.032%
94	15.725	2266	2267	2270	rBV3	1089	1111	0.28%	0.049%
95	15.804	2275	2280	2282	rBV6	1419	2316	0.58%	0.102%
96	15.828	2282	2284	2287	rBV4	1561	1198	0.30%	0.053%
97	16.035	2315	2318	2319	rBV3	970	1007	0.25%	0.044%
98	16.176	2339	2341	2344	rVB4	1245	1047	0.26%	0.046%
99	16.462	2385	2388	2389	rBV3	1159	1137	0.28%	0.050%
100	16.749	2433	2435	2439	rBV4	869	691	0.17%	0.030%

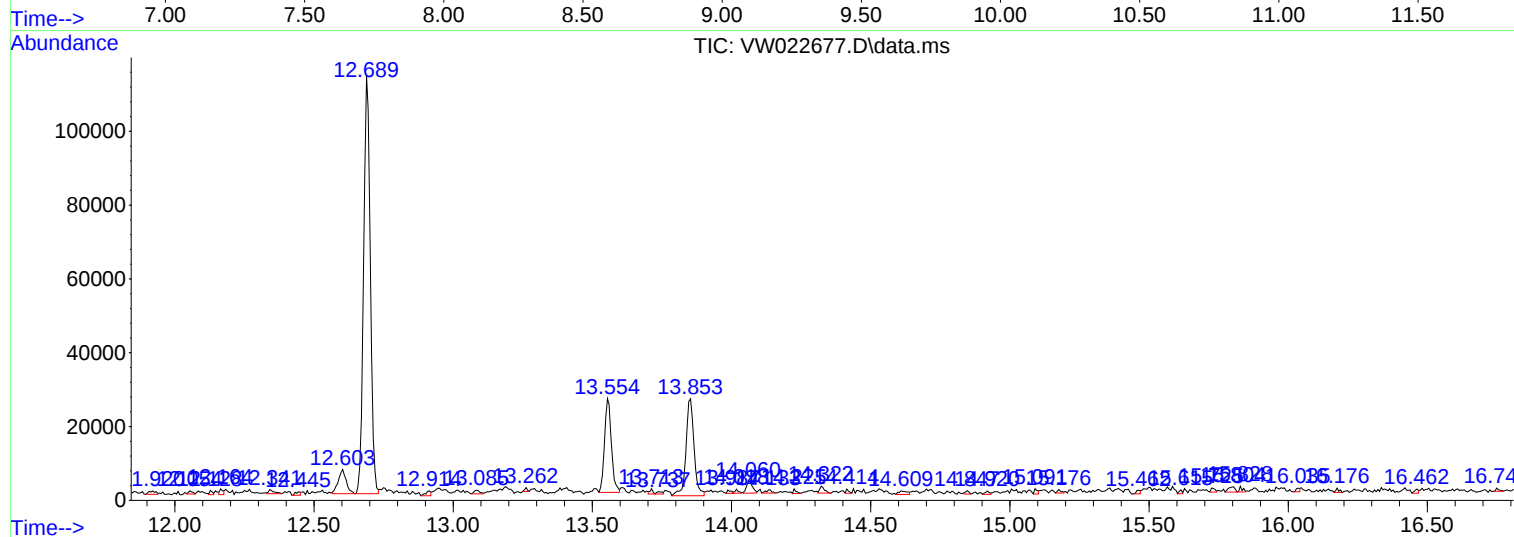
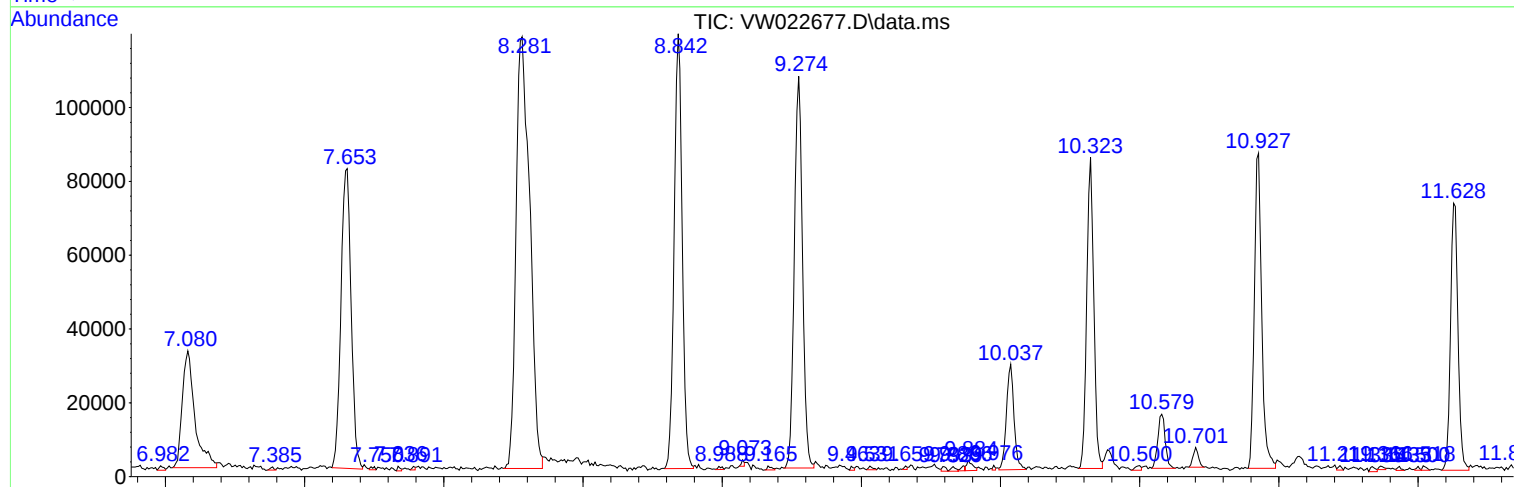
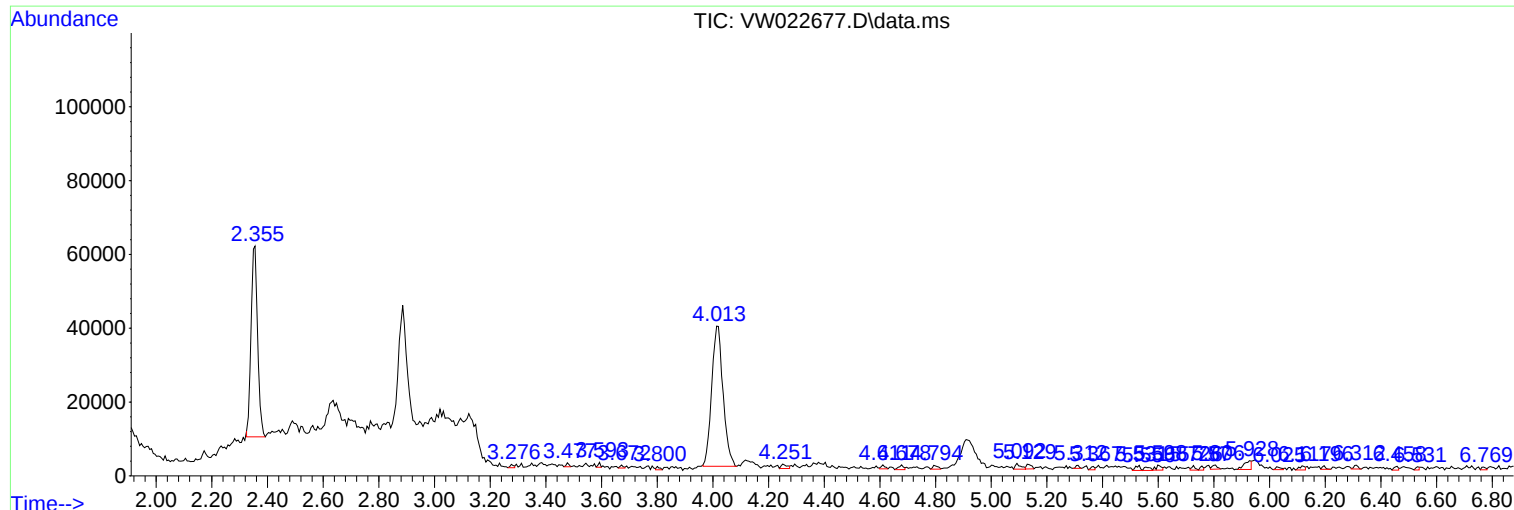
Sum of corrected areas: 2279011

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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM040722SMA.M
Quant Title : SFAM01.0

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



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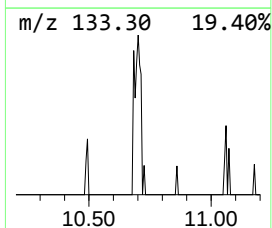
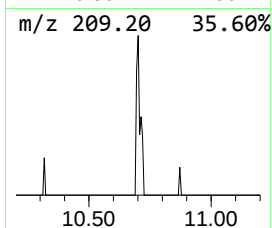
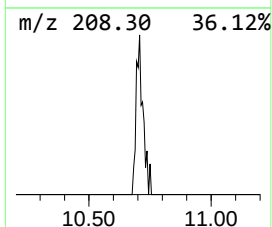
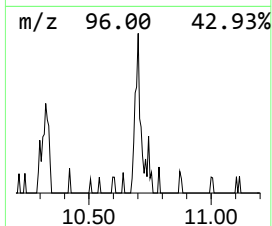
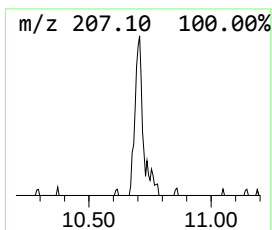
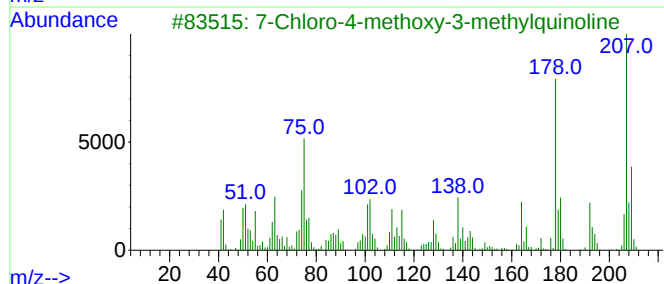
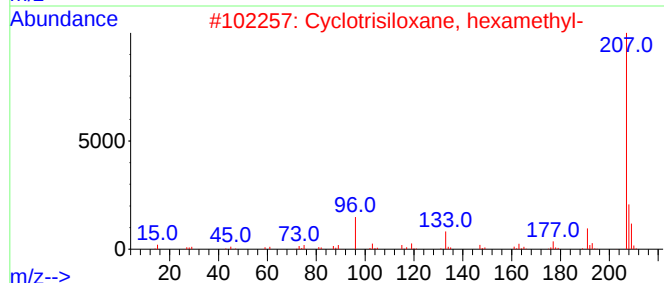
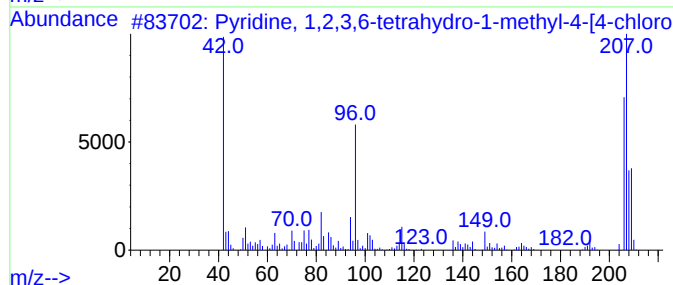
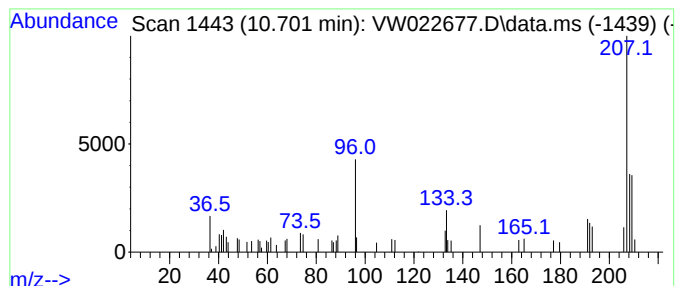
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM040722SMA.M
Quant Title : SFAM01.0

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

Peak Number 2 Pyridine, 1,2,3,6-tetrahydr... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.701	1.33 ug/L	7120	Chlorobenzene-d5	11.628

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pyridine, 1,2,3,6-tetrahydro-1-m...	207	C12H14ClN	005048-08-8	59
2			Cyclotrisiloxane, hexamethyl-	222	C6H18O3Si3	000541-05-9	50
3			7-Chloro-4-methoxy-3-methylquino...	207	C11H10ClNO	1000213-52-2	47
4			2-Ethylacridine	207	C15H13N	055751-83-2	43
5			2-Butenenitrile, 2-chloro-3-(4-m...	207	C11H10ClNO	1010305-66-7	43



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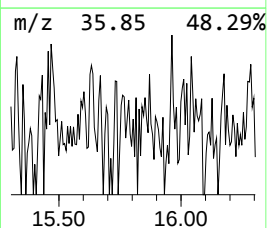
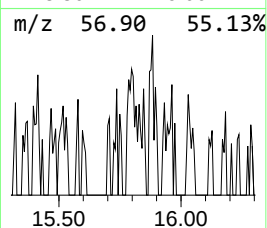
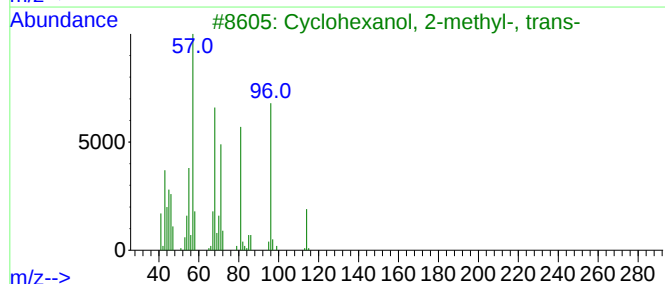
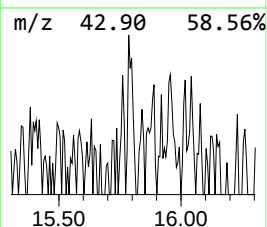
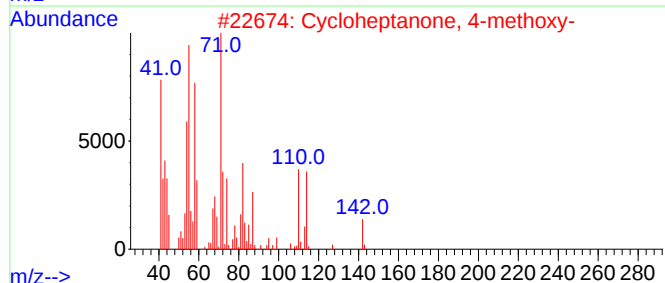
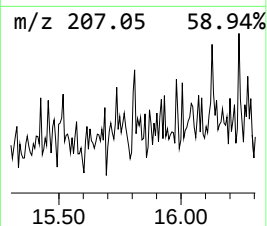
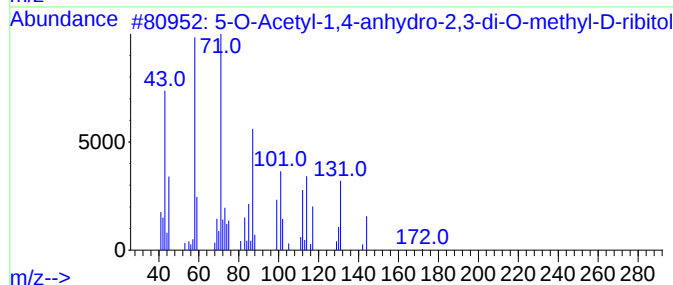
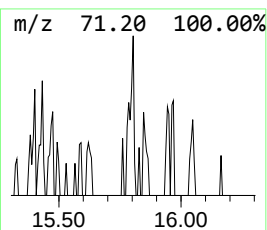
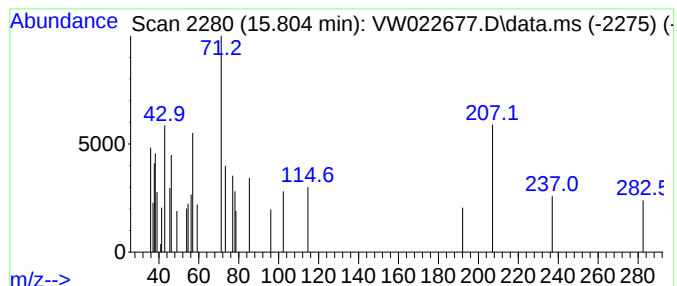
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Quant Title : SFAM01.0

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

Peak Number 5 unknown-01 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.804	1.34 ug/L	2316	1,4-Dichlorobenzene-d4	13.554

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	5-O-Acetyl-1,4-anhydro-2,3-di-O-...	204	C9H16O5	1000357-22-9	12		
2	Cycloheptanone, 4-methoxy-	142	C8H14O2	017429-01-5	10		
3	Cyclohexanol, 2-methyl-, trans-	114	C7H14O	007443-52-9	9		
4	1,3,5-Triazine, 2-chloro-4,6-bis...	207	C5H6ClN3S2	004407-40-3	9		
5	(Aminomethyl)cyclopropane	71	C4H9N	002516-47-4	7		



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TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Pyridine, 1,2,3...	10.701	1.3	ug/L	7120	2	11.628	134332	25.0
unknown-01	15.804	1.3	ug/L	2316	3	13.554	43280	25.0