

Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW080922\
 Data File : VW024211.D
 Acq On : 09 Aug 2022 16:01
 Operator : SY/VA
 Sample : N4110-04
 Misc : 5.26g/10mL/MSVOA_W/SOIL
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 C0ZF7

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\SFAMWLM080422SMA.M
 Title : SFAM01.0

Signal : TIC: VW024211.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	46	57	58	rBV7	12132	34834	3.70%	0.513%
2	2.337	66	71	77	rVB	133996	204106	21.68%	3.008%
3	2.483	92	95	104	rBV2	10569	20746	2.20%	0.306%
4	2.873	153	159	167	rVB	59146	119602	12.70%	1.762%
5	3.830	312	316	318	rBV3	617	818	0.09%	0.012%
6	3.891	324	326	329	rVB4	890	892	0.09%	0.013%
7	3.940	331	334	335	rBV3	1162	1319	0.14%	0.019%
8	4.001	335	344	356	rVV2	98851	289675	30.77%	4.269%
9	4.117	358	363	373	rVB2	13725	37223	3.95%	0.549%
10	4.348	400	401	403	rBV2	1050	1045	0.11%	0.015%
11	4.409	406	411	423	rVB2	6666	21787	2.31%	0.321%
12	4.757	465	468	469	rVB3	1054	887	0.09%	0.013%
13	4.781	469	472	474	rBV3	1048	1082	0.11%	0.016%
14	4.909	480	493	505	rBV3	22935	78265	8.31%	1.153%
15	4.989	505	506	510	rVB3	1356	1123	0.12%	0.017%
16	5.056	515	517	519	rBV3	1261	1031	0.11%	0.015%
17	5.123	519	528	533	rBV5	3774	13467	1.43%	0.198%
18	5.312	557	559	565	rVB6	1101	1330	0.14%	0.020%
19	5.379	565	570	571	rBV4	962	1146	0.12%	0.017%
20	5.403	572	574	577	rVV4	1013	997	0.11%	0.015%
21	5.440	577	580	584	rVB4	1565	2091	0.22%	0.031%
22	5.592	604	605	608	rVB2	1294	809	0.09%	0.012%
23	5.860	644	649	650	rBV4	800	1019	0.11%	0.015%
24	5.934	650	661	669	rVV6	8926	28023	2.98%	0.413%
25	6.031	675	677	679	rBV2	1184	1294	0.14%	0.019%
26	6.208	705	706	709	rVB2	1238	903	0.10%	0.013%
27	6.263	712	715	717	rVB3	1358	1180	0.13%	0.017%
28	6.287	717	719	721	rBV2	970	968	0.10%	0.014%
29	6.366	731	732	735	rBV2	1171	1226	0.13%	0.018%
30	6.702	782	787	789	rVB5	747	1116	0.12%	0.016%
31	6.750	792	795	798	rVB4	755	1028	0.11%	0.015%
32	6.842	807	810	812	rBV3	740	886	0.09%	0.013%
33	6.891	816	818	820	rBV3	957	976	0.10%	0.014%
34	6.915	820	822	826	rVB4	756	883	0.09%	0.013%
35	6.976	830	832	835	rVB3	1097	910	0.10%	0.013%
36	7.080	840	849	855	rBV	32286	89226	9.48%	1.315%

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Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM080422SMA.M
 Title : SFAM01.0

37	7.153	855	861	873	rVB	48341	135090	14.35%	1.991%
38	7.464	907	912	918	rBV6	902	1913	0.20%	0.028%
39	7.519	918	921	924	rBV4	1202	2048	0.22%	0.030%
40	7.543	924	925	928	rVB3	1489	982	0.10%	0.014%
41	7.580	928	931	932	rBV3	937	851	0.09%	0.013%
42	7.647	932	942	952	rVV	149354	364660	38.73%	5.374%
43	7.964	992	994	996	rVB3	1375	1081	0.11%	0.016%
44	8.177	1026	1029	1031	rVB3	721	880	0.09%	0.013%
45	8.275	1031	1045	1059	rBV2	293280	941531	100.00%	13.874%
46	8.543	1088	1089	1092	rBV3	1315	1019	0.11%	0.015%
47	8.701	1113	1115	1118	rBV4	1059	1154	0.12%	0.017%
48	8.842	1129	1138	1148	rBV	226732	455424	48.37%	6.711%
49	8.951	1153	1156	1158	rBV4	1336	1538	0.16%	0.023%
50	9.037	1166	1170	1173	rBV5	2602	3687	0.39%	0.054%
51	9.274	1201	1209	1220	rVV	199850	404007	42.91%	5.953%
52	9.402	1228	1230	1234	rVB5	1372	1820	0.19%	0.027%
53	9.445	1234	1237	1239	rBV4	976	970	0.10%	0.014%
54	9.506	1245	1247	1248	rBV2	1099	746	0.08%	0.011%
55	9.677	1272	1275	1278	rVB5	1116	1472	0.16%	0.022%
56	9.707	1278	1280	1284	rBV5	872	1100	0.12%	0.016%
57	9.841	1299	1302	1303	rBV2	923	862	0.09%	0.013%
58	9.878	1305	1308	1309	rBV3	1360	1225	0.13%	0.018%
59	9.939	1316	1318	1319	rBV2	843	840	0.09%	0.012%
60	9.975	1322	1324	1327	rBV3	1372	1951	0.21%	0.029%
61	10.036	1327	1334	1342	rVB	92274	165973	17.63%	2.446%
62	10.323	1373	1381	1390	rBV	362494	656947	69.77%	9.681%
63	10.475	1404	1406	1407	rBV2	1121	914	0.10%	0.013%
64	10.579	1416	1423	1431	rBV	63029	114781	12.19%	1.691%
65	10.701	1438	1443	1451	rVB3	17134	31852	3.38%	0.469%
66	10.762	1451	1453	1456	rBV4	799	878	0.09%	0.013%
67	10.841	1462	1466	1471	rBV7	2536	4681	0.50%	0.069%
68	10.927	1473	1480	1490	rBV	130353	226369	24.04%	3.336%
69	11.067	1497	1503	1510	rVB2	22073	39798	4.23%	0.586%
70	11.176	1515	1521	1526	rVB8	4985	9325	0.99%	0.137%
71	11.341	1546	1548	1550	rBV3	1082	1155	0.12%	0.017%
72	11.433	1554	1563	1569	rVB4	13614	31485	3.34%	0.464%
73	11.506	1573	1575	1580	rVB5	1512	1744	0.19%	0.026%
74	11.573	1582	1586	1588	rBV4	854	1296	0.14%	0.019%
75	11.628	1588	1595	1605	rBV	307018	529292	56.22%	7.800%
76	11.847	1625	1631	1632	rBV6	1809	3018	0.32%	0.044%

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Peak separation: 5

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 Title : SFAM01.0

77	11.945	1644	1647	1648	rVB3	1245	915	0.10%	0.013%
78	11.999	1654	1656	1659	rBV4	1141	1582	0.17%	0.023%
79	12.134	1673	1678	1683	rBV8	7504	13309	1.41%	0.196%
80	12.249	1692	1697	1702	rVB3	14498	25569	2.72%	0.377%
81	12.335	1702	1711	1712	rBV4	8685	18534	1.97%	0.273%
82	12.359	1712	1715	1722	rVB3	12366	19950	2.12%	0.294%
83	12.414	1722	1724	1726	rBV3	1653	1095	0.12%	0.016%
84	12.499	1727	1738	1742	rBV9	12377	36375	3.86%	0.536%
85	12.603	1750	1755	1761	rVB2	27563	49698	5.28%	0.732%
86	12.688	1763	1769	1777	rBV	191272	321839	34.18%	4.743%
87	12.823	1790	1791	1792	rBV	1952	1368	0.15%	0.020%
88	12.865	1793	1798	1801	rVV7	6309	11957	1.27%	0.176%
89	12.932	1803	1809	1815	rBV6	22226	56575	6.01%	0.834%
90	13.072	1829	1832	1836	rBV3	17915	25253	2.68%	0.372%
91	13.121	1837	1840	1843	rBV4	18105	26324	2.80%	0.388%
92	13.164	1843	1847	1852	rVB2	38596	61347	6.52%	0.904%
93	13.292	1864	1868	1872	rVB2	23219	32631	3.47%	0.481%
94	13.554	1906	1911	1920	rVB	269446	430356	45.71%	6.342%
95	13.853	1953	1960	1967	rBV	261932	502995	53.42%	7.412%
96	15.164	2173	2175	2176	rBV2	2688	1869	0.20%	0.028%
97	15.322	2198	2201	2202	rBV3	2335	2004	0.21%	0.030%
98	15.432	2214	2219	2226	rVB2	20879	32589	3.46%	0.480%
99	15.755	2270	2272	2273	rBV2	1904	953	0.10%	0.014%
100	16.084	2324	2326	2328	rBV3	1197	844	0.09%	0.012%

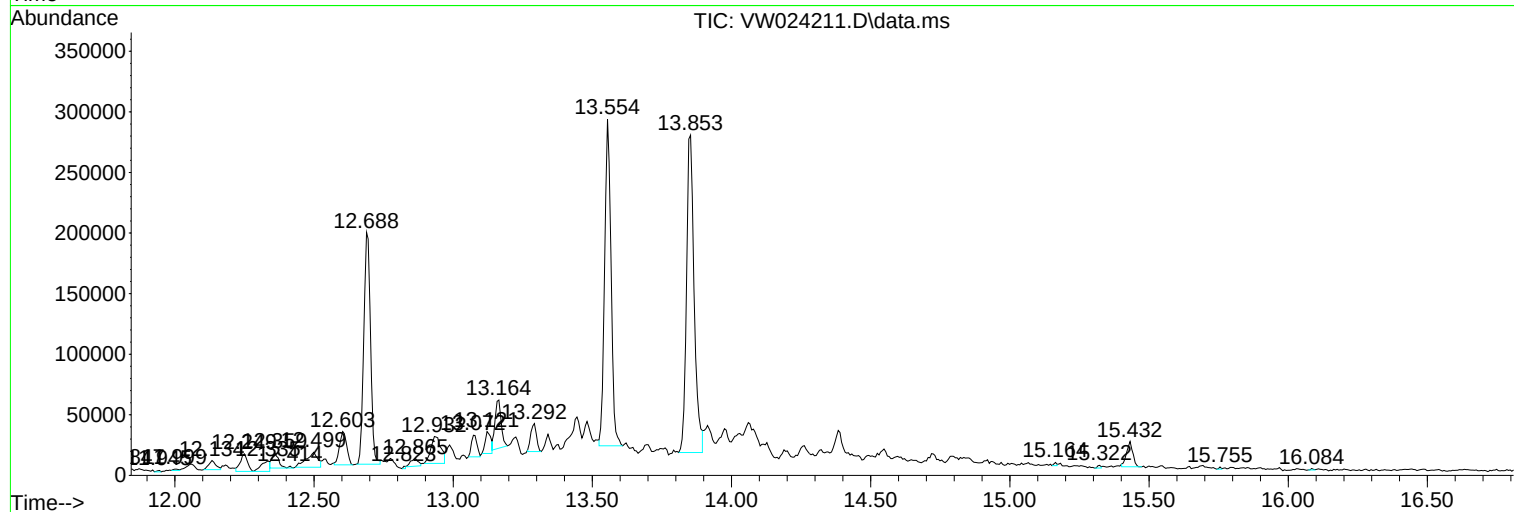
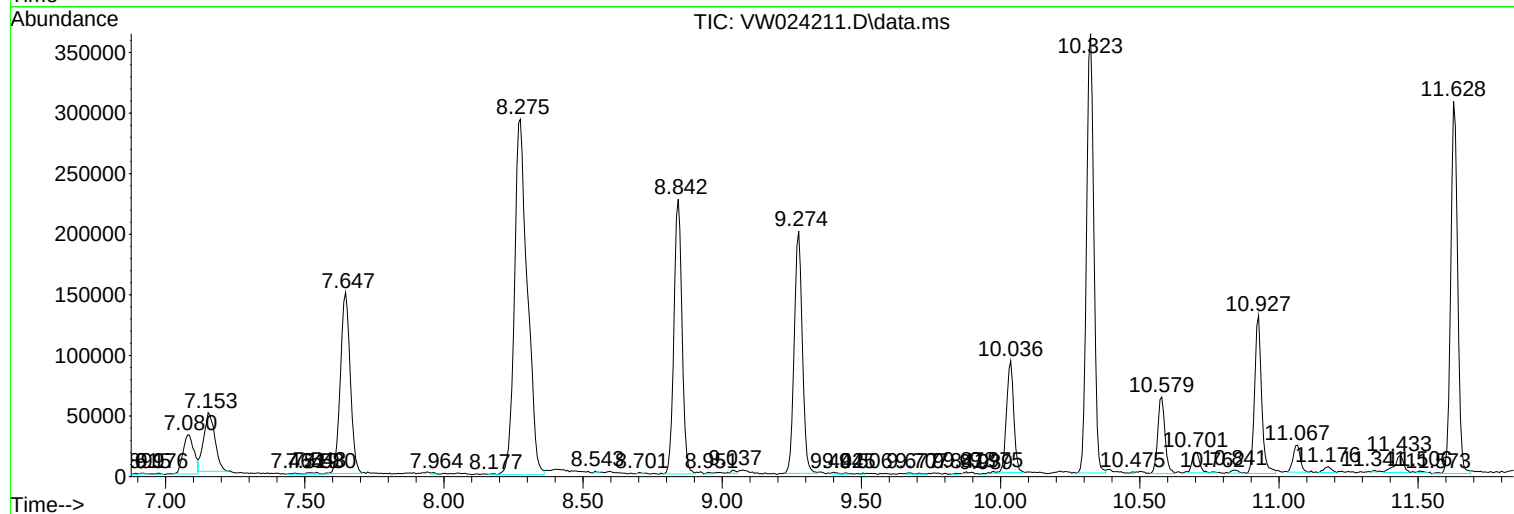
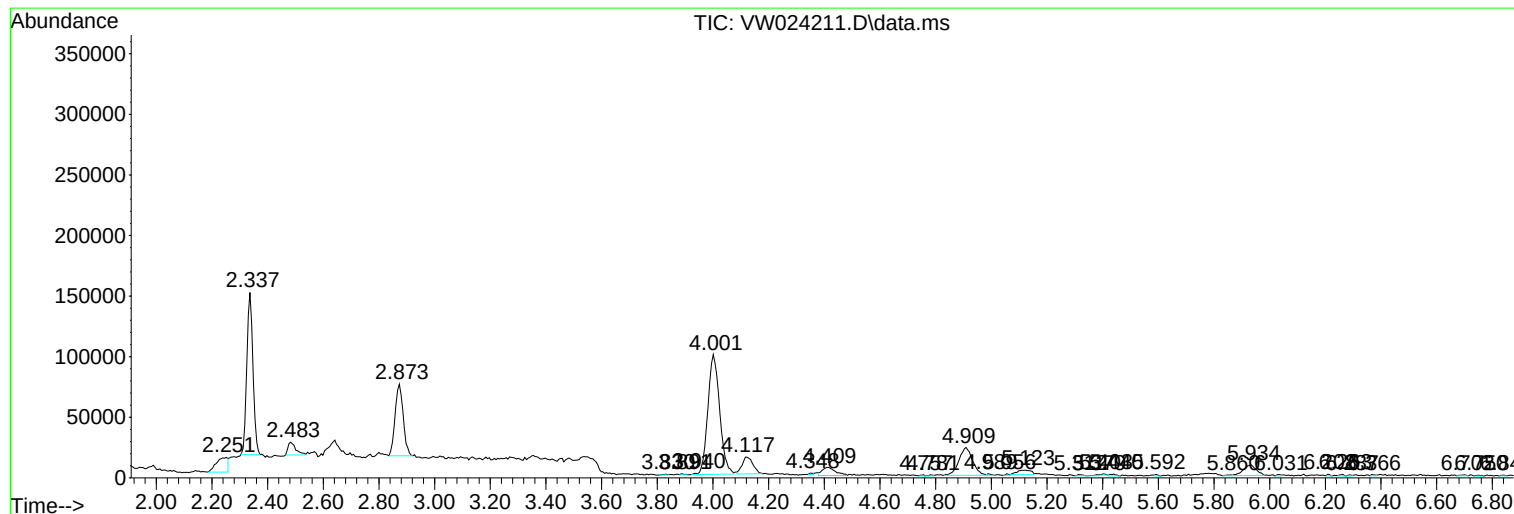
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ALS Vial : 1 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM080422SMA.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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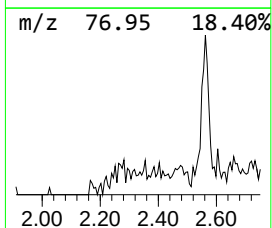
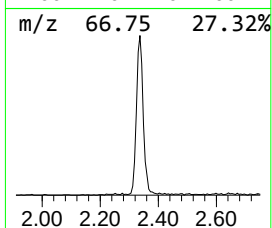
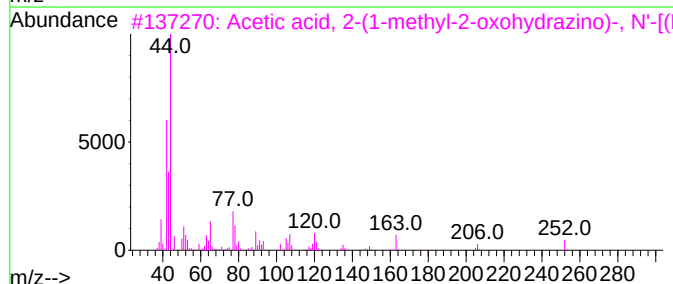
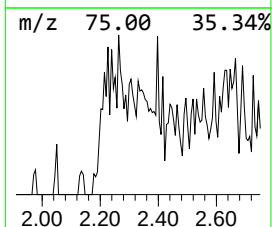
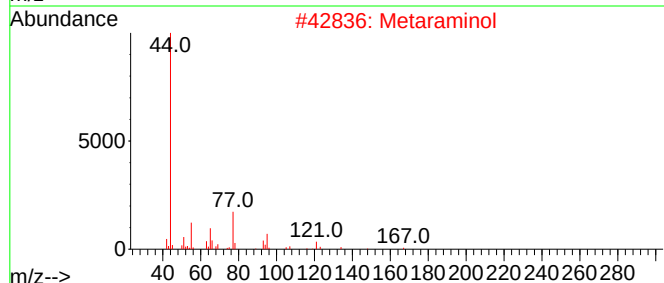
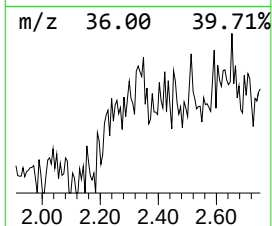
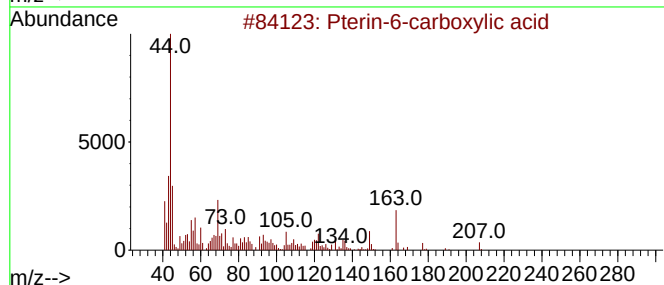
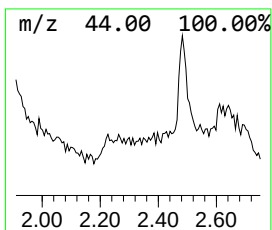
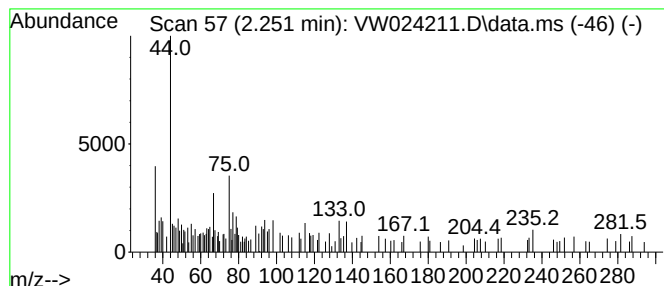
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM080422SMA.M
Quant Title : SFAM01.0

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

Peak Number 1 Pterin-6-carboxylic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.251	1.91 ug/L	34834	1,4-Difluorobenzene	8.842

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pterin-6-carboxylic acid	207	C7H5N5O3	000948-60-7	52
2			Metaraminol	167	C9H13NO2	000054-49-9	43
3			Acetic acid, 2-(1-methyl-2-oxohy...	252	C10H12N4O4	1000327-59-6	43
4			Pyrazole[4,5-b]imidazole, 1-form...	296	C12H16N4O5	1000127-72-0	40
5			Pyridine-3-carboxamide, 1,2-dihy...	182	C8H10N2O5	079927-21-2	40



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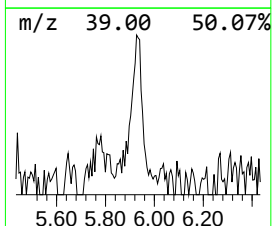
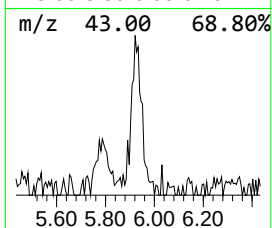
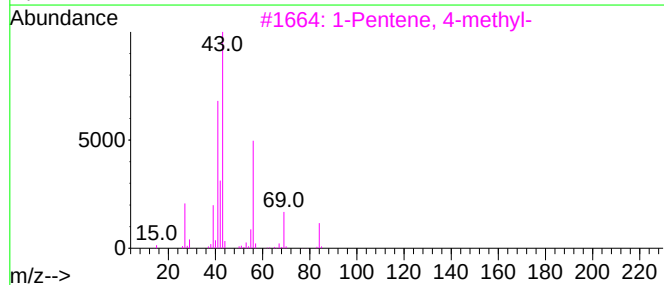
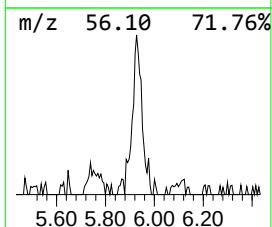
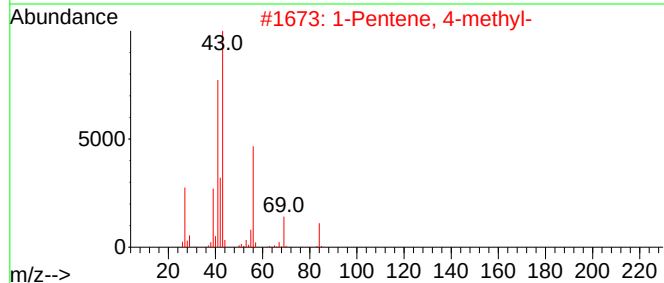
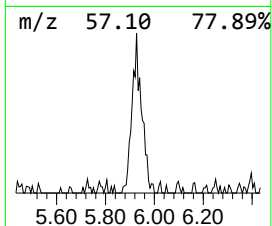
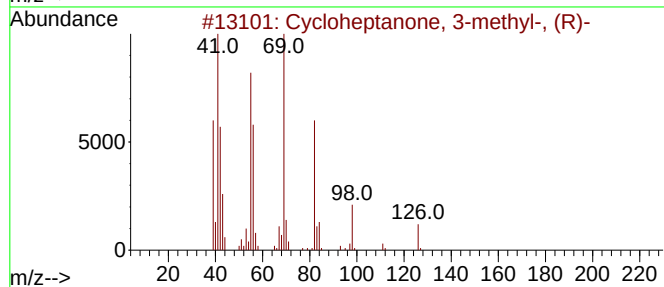
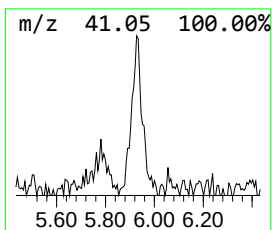
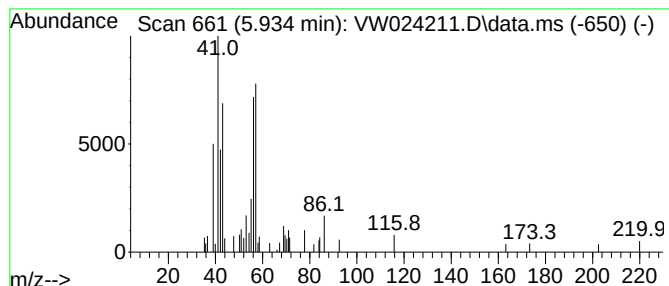
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM080422SMA.M
Quant Title : SFAM01.0

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

Peak Number 2 unknown-01 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.934	1.54 ug/L	28023	1,4-Difluorobenzene	8.842

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cycloheptanone, 3-methyl-, (R)-	126	C8H14O	013609-58-0	45
2			1-Pentene, 4-methyl-	84	C6H12	000691-37-2	43
3			1-Pentene, 4-methyl-	84	C6H12	000691-37-2	43
4			3-Buten-1-ol, 3-methyl-	86	C5H10O	000763-32-6	43
5			1-Heptanethiol	132	C7H16S	001639-09-4	42



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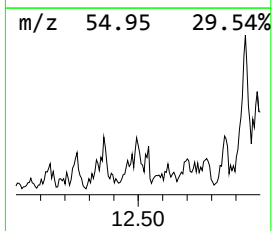
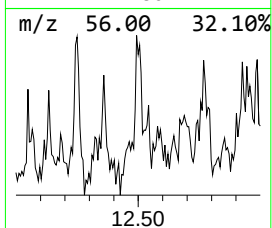
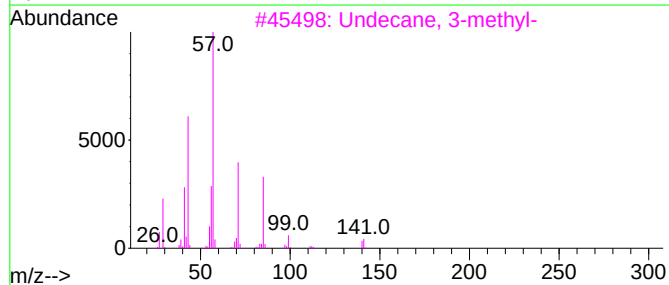
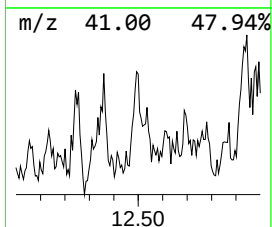
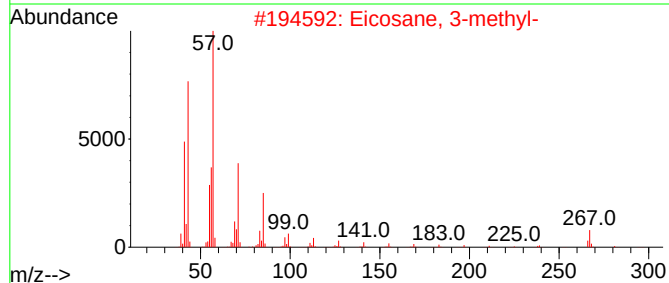
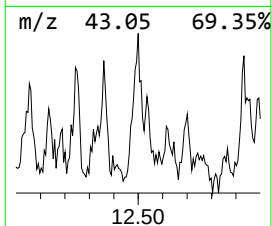
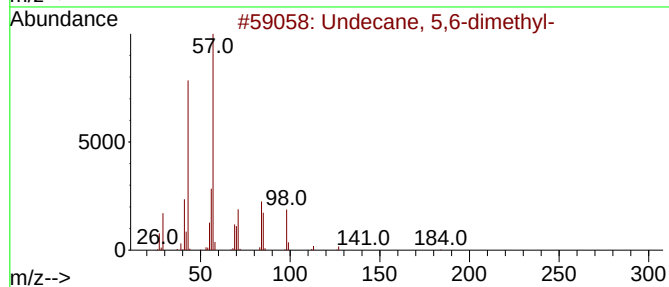
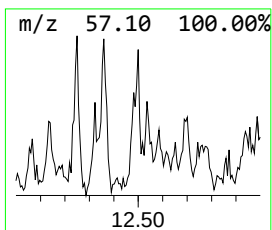
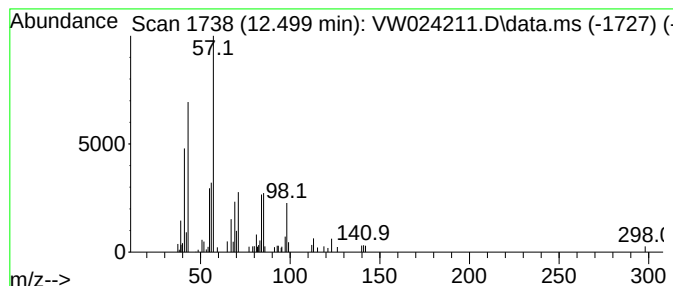
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM080422SMA.M
Quant Title : SFAM01.0

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

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.499	1.72 ug/L	36375	Chlorobenzene-d5	11.628

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Undecane, 5,6-dimethyl-	184	C13H28	017615-91-7	59
2			Eicosane, 3-methyl-	296	C21H44	006418-46-8	47
3			Undecane, 3-methyl-	170	C12H26	001002-43-3	47
4			Oxalic acid, isobutyl tetradecyl...	342	C20H38O4	1000309-37-9	47
5			Octane, 4-ethyl-	142	C10H22	015869-86-0	43



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW080922\
Data File : VW024211.D
Acq On : 09 Aug 2022 16:01
Operator : SY/VA
Sample : N4110-04
Misc : 5.26g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
C0ZF7

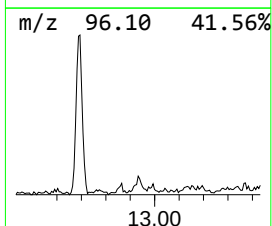
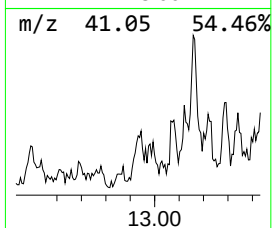
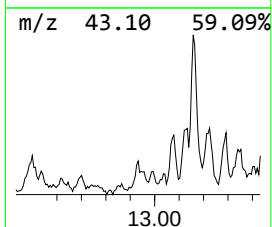
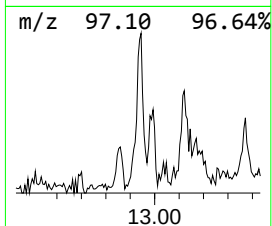
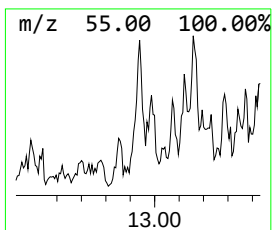
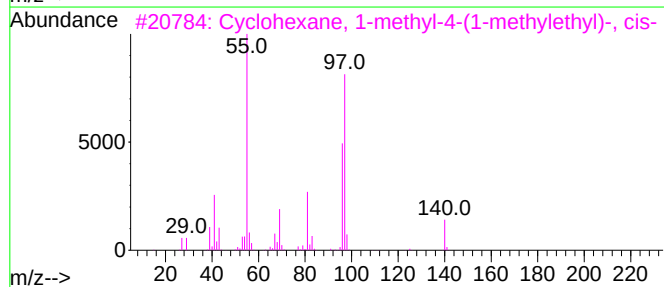
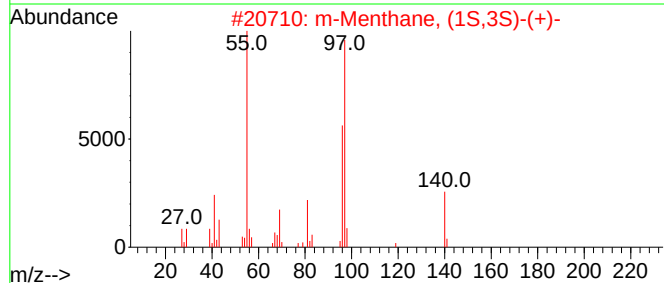
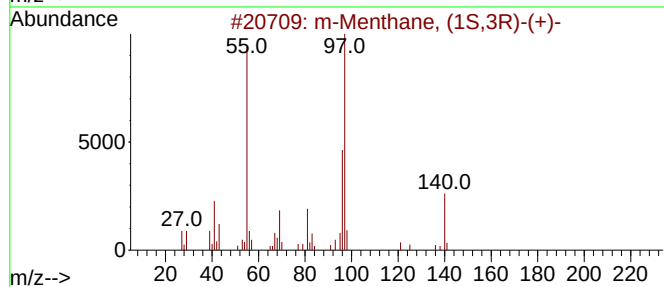
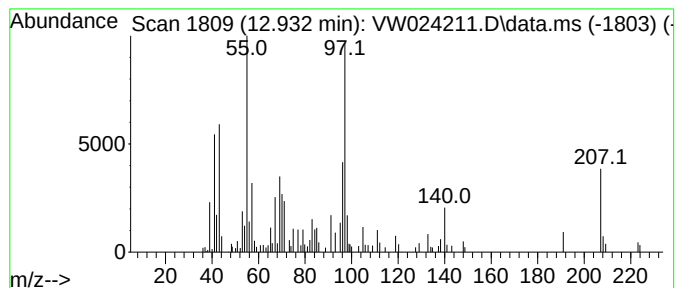
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Quant Title : SFAM01.0

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

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.932	3.29 ug/L	56575	1,4-Dichlorobenzene-d4	13.554

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			m-Menthane, (1S,3R)-(+)-	140	C10H20	013837-66-6	62
2			m-Menthane, (1S,3S)-(+)-	140	C10H20	013837-67-7	52
3			Cyclohexane, 1-methyl-4-(1-methy...	140	C10H20	006069-98-3	50
4			Oxalic acid, cyclohexylmethyl oc...	298	C17H30O4	1000309-68-4	49
5			Cyclohexane, 1-(cyclohexylmethyl...	194	C14H26	054823-95-9	47



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW080922\
Data File : VW024211.D
Acq On : 09 Aug 2022 16:01
Operator : SY/VA
Sample : N4110-04
Misc : 5.26g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
C0ZF7

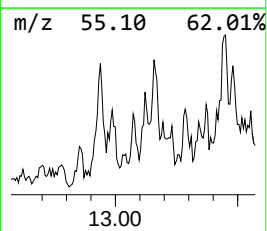
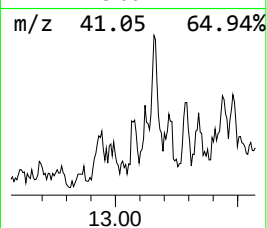
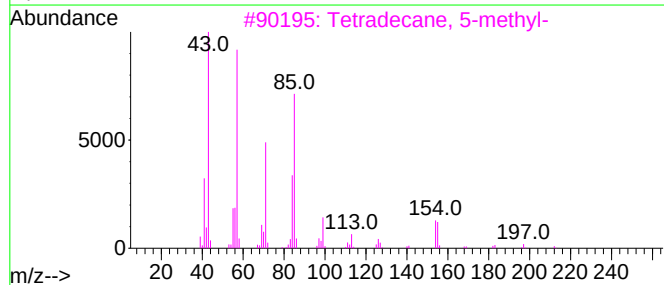
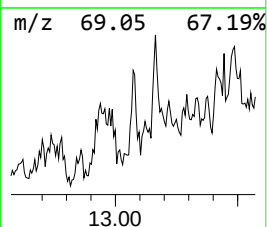
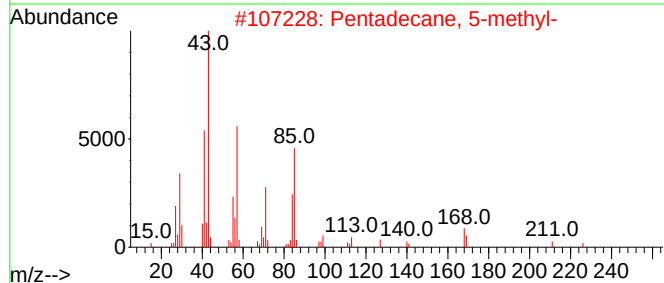
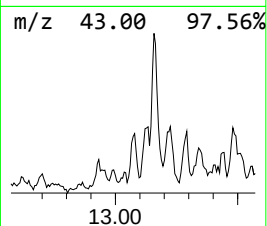
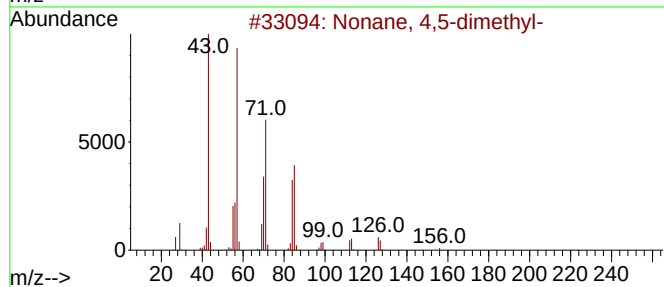
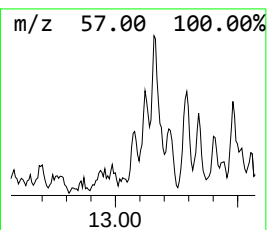
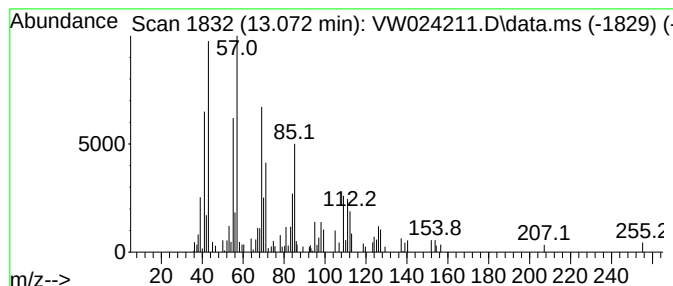
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Quant Title : SFAM01.0

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

Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.073	1.47 ug/L	25253	1,4-Dichlorobenzene-d4	13.554

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonane, 4,5-dimethyl-	156	C11H24	017302-23-7	53
2			Pentadecane, 5-methyl-	226	C16H34	025117-33-3	47
3			Tetradecane, 5-methyl-	212	C15H32	025117-32-2	47
4			Pentadecane	212	C15H32	000629-62-9	43
5			Sulfurous acid, butyl decyl ester	278	C14H30O3S	1000309-17-7	38



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW080922\
Data File : VW024211.D
Acq On : 09 Aug 2022 16:01
Operator : SY/VA
Sample : N4110-04
Misc : 5.26g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
C0ZF7

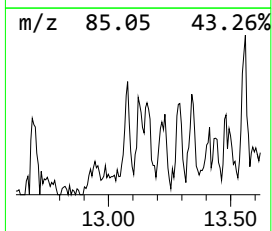
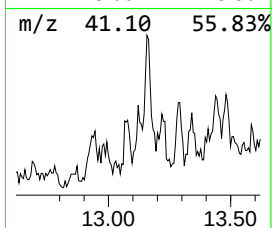
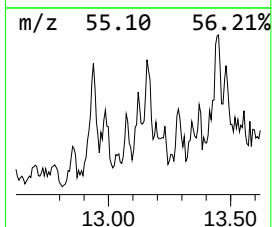
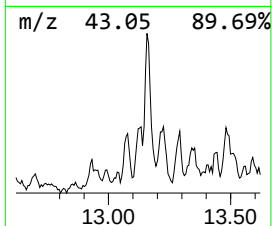
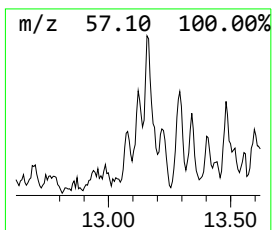
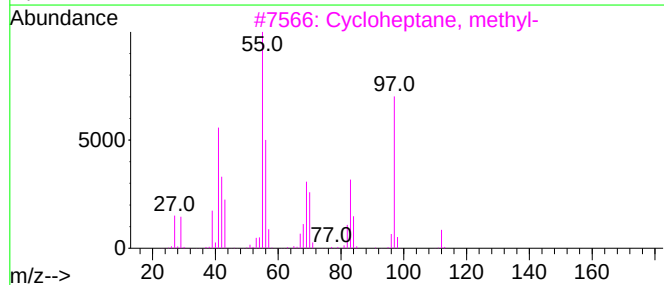
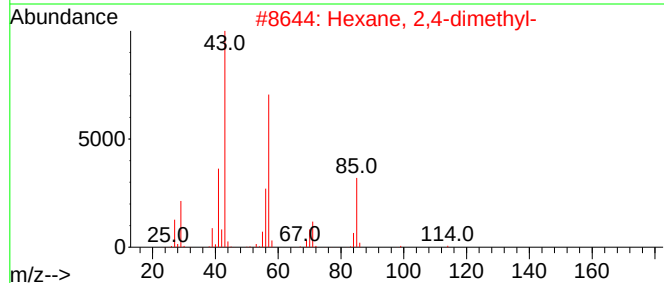
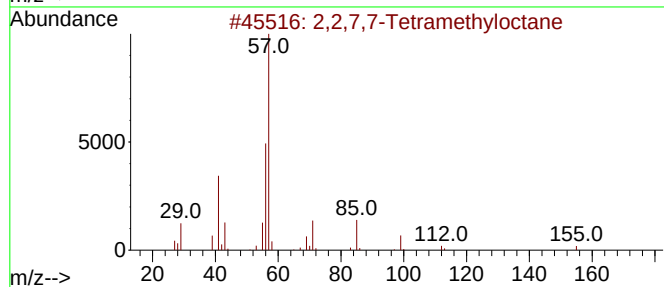
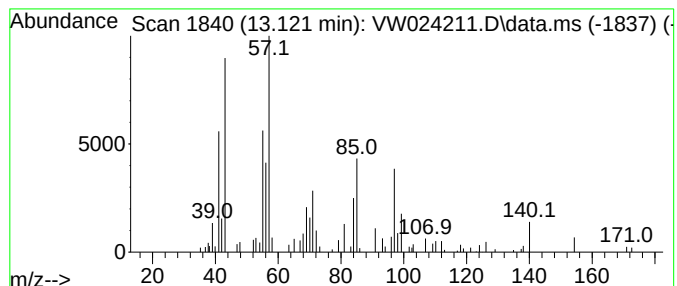
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Quant Title : SFAM01.0

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

Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.121	1.53 ug/L	26324	1,4-Dichlorobenzene-d4	13.554

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2,2,7,7-Tetramethyloctane	170	C12H26	001071-31-4	47
2			Hexane, 2,4-dimethyl-	114	C8H18	000589-43-5	43
3			Cycloheptane, methyl-	112	C8H16	004126-78-7	42
4			Ether, heptyl hexyl	200	C13H28O	007289-40-9	38
5			4-Decene, 7-methyl-, (E)-	154	C11H22	062338-48-1	30



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW080922\
Data File : VW024211.D
Acq On : 09 Aug 2022 16:01
Operator : SY/VA
Sample : N4110-04
Misc : 5.26g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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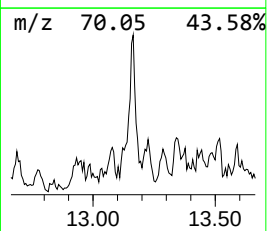
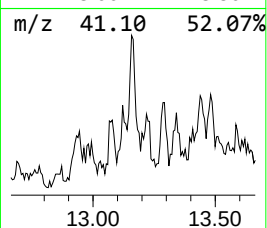
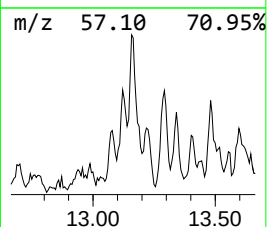
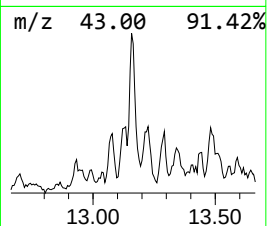
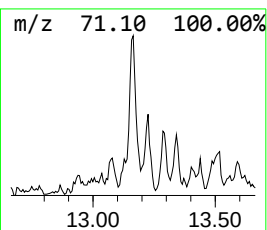
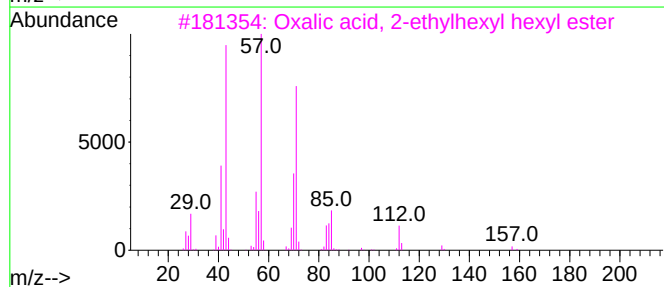
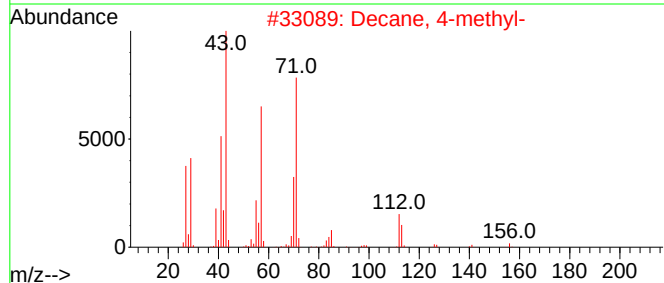
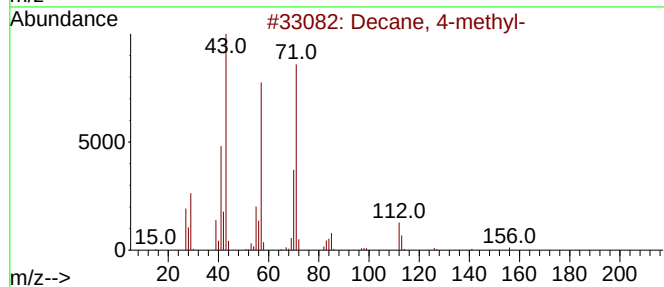
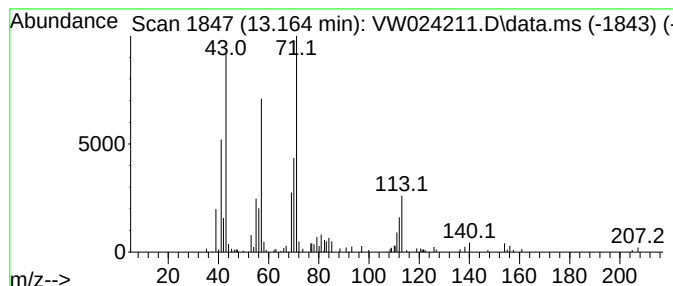
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Quant Title : SFAM01.0

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

Peak Number 12 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.164	3.56 ug/L	61347	1,4-Dichlorobenzene-d4	13.554

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Decane, 4-methyl-	156	C11H24	002847-72-5	68
2			Decane, 4-methyl-	156	C11H24	002847-72-5	59
3			Oxalic acid, 2-ethylhexyl hexyl ...	286	C16H30O4	1000309-38-9	59
4			4-Methyl-3-heptanol, chlorodiflu...	242	C10H17ClF2O2	1000447-28-3	53
5			Oxalic acid, 2-ethylhexyl pentyl...	272	C15H28O4	1000309-38-7	53



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW080922\
Data File : VW024211.D
Acq On : 09 Aug 2022 16:01
Operator : SY/VA
Sample : N4110-04
Misc : 5.26g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
C0ZF7

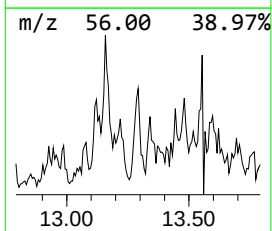
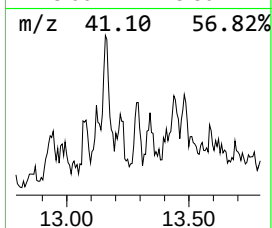
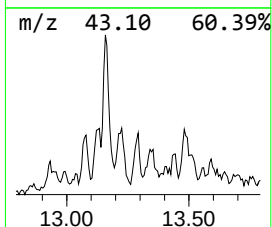
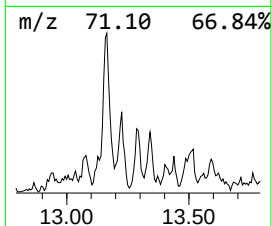
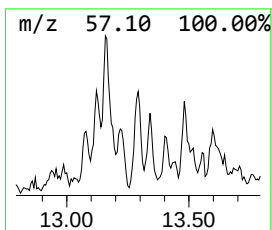
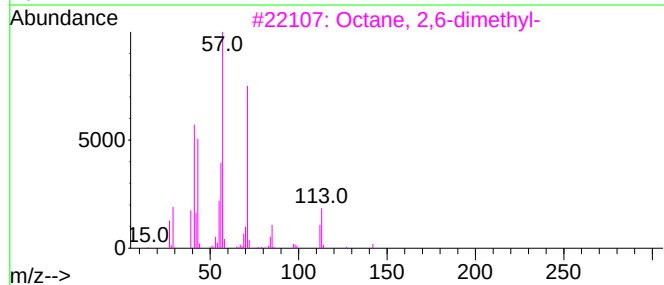
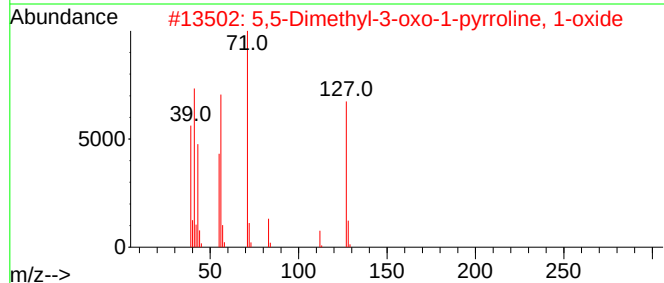
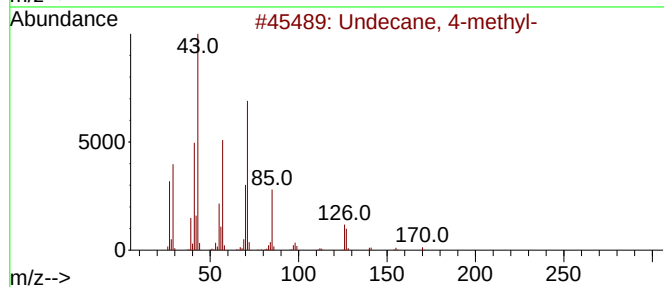
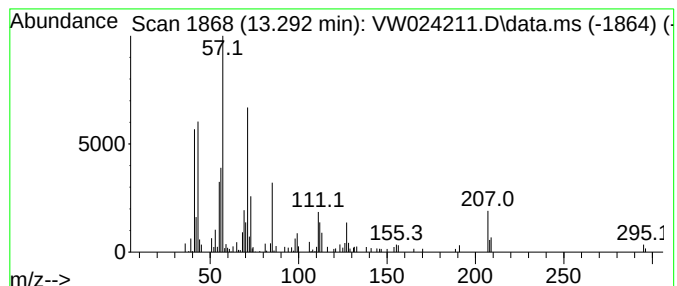
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Quant Title : SFAM01.0

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

Peak Number 13 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.292	1.90 ug/L	32631	1,4-Dichlorobenzene-d4	13.554

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Undecane, 4-methyl-	170	C12H26	002980-69-0	38
2			5,5-Dimethyl-3-oxo-1-pyrroline, ...	127	C6H9NO2	1000305-98-3	38
3			Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	38
4			Butane, 2,2-dimethyl-	86	C6H14	000075-83-2	35
5			Sulfurous acid, nonyl 2-propyl e...	250	C12H26O3S	1000309-12-0	35



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW080922\
Data File : VW024211.D
Acq On : 09 Aug 2022 16:01
Operator : SY/VA
Sample : N4110-04
Misc : 5.26g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
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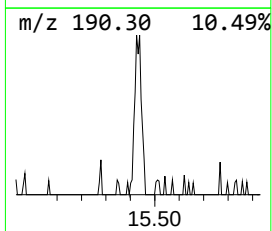
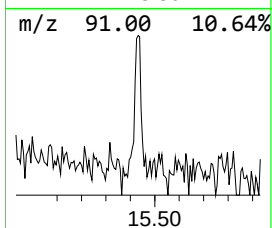
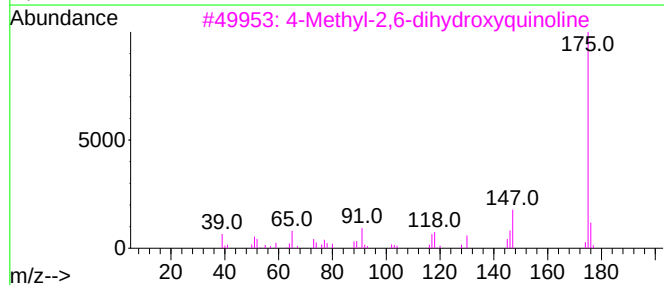
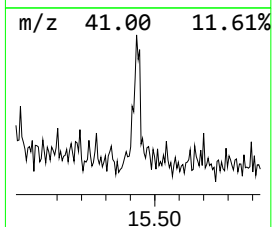
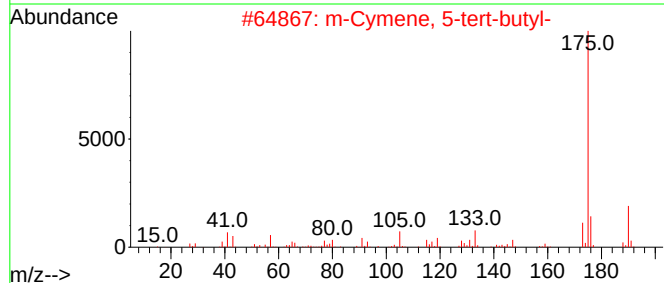
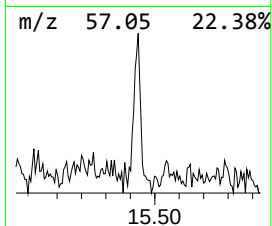
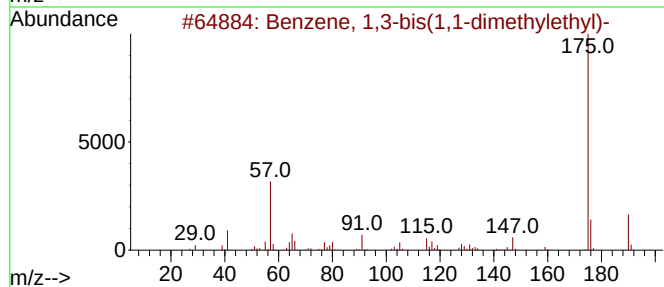
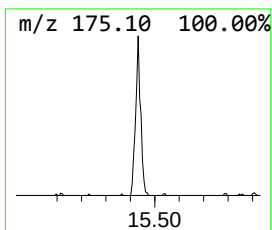
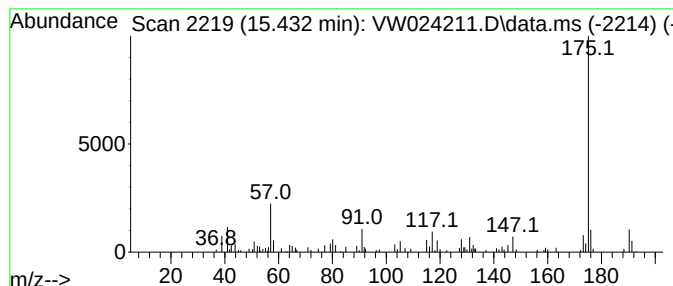
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Quant Title : SFAM01.0

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

Peak Number 14 Benzene, 1,3-bis(1,1-dimeth... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.432	1.89 ug/L	32589	1,4-Dichlorobenzene-d4	13.554

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,3-bis(1,1-dimethyleth...	190	C14H22	001014-60-4	87
2			m-Cymene, 5-tert-butyl-	190	C14H22	029577-19-3	80
3			4-Methyl-2,6-dihydroxyquinoline	175	C10H9NO2	034982-01-9	68
4			4-Methyl-2,6-dihydroxyquinoline	175	C10H9NO2	034982-01-9	64
5			Benzene, 1,3-bis(1,1-dimethyleth...	190	C14H22	001014-60-4	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW080922\
Data File : VW024211.D
Acq On : 09 Aug 2022 16:01
Operator : SY/VA
Sample : N4110-04
Misc : 5.26g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
C0ZF7

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM080422SMA.M
Quant Title : SFAM01.0

TIC Library : C:\Database\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
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unknown-01	5.934	1.5	ug/L	28023	1	8.842	455424	25.0
(DEL) Alkane: S...	12.499	1.7	ug/L	36375	2	11.628	529292	25.0
(DEL) Alkane: S...	12.932	3.3	ug/L	56575	3	13.554	430356	25.0
(DEL) Alkane: S...	13.073	1.5	ug/L	25253	3	13.554	430356	25.0
(DEL) Alkane: S...	13.121	1.5	ug/L	26324	3	13.554	430356	25.0
(DEL) Alkane: S...	13.164	3.6	ug/L	61347	3	13.554	430356	25.0
(DEL) Alkane: S...	13.292	1.9	ug/L	32631	3	13.554	430356	25.0
Benzene, 1,3-bi...	15.432	1.9	ug/L	32589	3	13.554	430356	25.0