

Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW051222\
 Data File : VW023011.D
 Acq On : 12 May 2022 19:06
 Operator : SY/VA
 Sample : N2767-09
 Misc : 4.76g/10mL/MSVOA_W/SOIL
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 BGPM4

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\SFAMWLM050222SMA.M
 Title : SFAM01.0

Signal : TIC: VW023011.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.221	48	52	54	rBV3	4283	5208	0.36%	0.051%
2	2.349	68	73	82	rBV	93448	160363	11.19%	1.573%
3	2.885	155	161	170	rVB	68755	150844	10.53%	1.480%
4	3.245	219	220	223	rVB3	1512	1305	0.09%	0.013%
5	3.385	239	243	248	rVV6	1963	4294	0.30%	0.042%
6	3.495	258	261	264	rBV4	684	1205	0.08%	0.012%
7	3.611	278	280	282	rVB3	1096	913	0.06%	0.009%
8	3.745	297	302	304	rBV5	933	1368	0.10%	0.013%
9	3.769	304	306	310	rVB5	880	907	0.06%	0.009%
10	4.013	335	346	360	rBV	97025	307491	21.46%	3.017%
11	4.190	371	375	376	rBV3	1229	1912	0.13%	0.019%
12	4.239	381	383	386	rVB4	1021	1079	0.08%	0.011%
13	4.397	402	409	418	rBV2	6338	18386	1.28%	0.180%
14	4.915	481	494	504	rBV2	19571	73432	5.12%	0.720%
15	5.037	511	514	517	rVB3	1199	1528	0.11%	0.015%
16	5.184	536	538	541	rVV3	1417	1496	0.10%	0.015%
17	5.940	653	662	673	rVB3	10458	31513	2.20%	0.309%
18	6.068	681	683	684	rBV	983	964	0.07%	0.009%
19	6.086	684	686	688	rVV	1121	1012	0.07%	0.010%
20	6.202	703	705	707	rBV3	754	1001	0.07%	0.010%
21	6.281	716	718	722	rVB3	1013	1261	0.09%	0.012%
22	6.427	739	742	744	rVB4	839	903	0.06%	0.009%
23	6.482	747	751	756	rBV7	1131	2582	0.18%	0.025%
24	6.622	773	774	779	rVB3	964	973	0.07%	0.010%
25	6.683	783	784	787	rVB3	1460	1001	0.07%	0.010%
26	6.824	803	807	811	rBV3	760	1715	0.12%	0.017%
27	6.891	815	818	822	rBV4	869	1348	0.09%	0.013%
28	6.976	830	832	837	rBV5	967	1833	0.13%	0.018%
29	7.074	841	848	854	rVV	43836	119925	8.37%	1.177%
30	7.141	855	859	872	rVV2	42236	115358	8.05%	1.132%
31	7.647	932	942	957	rVV	195843	492958	34.40%	4.836%
32	7.756	957	960	962	rVB4	1016	1188	0.08%	0.012%
33	7.951	988	992	996	rVB5	2337	4200	0.29%	0.041%
34	8.006	996	1001	1002	rBV5	985	1167	0.08%	0.011%
35	8.201	1030	1033	1034	rBV3	980	962	0.07%	0.009%
36	8.274	1034	1045	1060	rVV2	347446	1084089	75.65%	10.635%

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Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM050222SMA.M
 Title : SFAM01.0

37	8.628	1100	1103	1105	rBV4	1053	1239	0.09%	0.012%
38	8.689	1108	1113	1117	rVV5	1530	2101	0.15%	0.021%
39	8.841	1127	1138	1150	rBV	327093	658789	45.97%	6.463%
40	9.030	1166	1169	1172	rBV3	1707	2681	0.19%	0.026%
41	9.274	1199	1209	1216	rBV	255089	520625	36.33%	5.108%
42	9.457	1236	1239	1242	rVB4	1132	1345	0.09%	0.013%
43	9.524	1248	1250	1251	rVB2	1491	929	0.06%	0.009%
44	9.573	1254	1258	1259	rBV4	1370	1921	0.13%	0.019%
45	9.671	1268	1274	1277	rBV6	1416	2730	0.19%	0.027%
46	9.762	1283	1289	1290	rBV5	1824	3037	0.21%	0.030%
47	9.890	1305	1310	1316	rVV6	5327	10484	0.73%	0.103%
48	9.957	1316	1321	1324	rVV2	8833	14973	1.04%	0.147%
49	10.036	1327	1334	1341	rVV	119252	222832	15.55%	2.186%
50	10.097	1341	1344	1356	rVB4	11017	23640	1.65%	0.232%
51	10.213	1358	1363	1366	rBV5	4046	7938	0.55%	0.078%
52	10.323	1373	1381	1388	rBV	455715	824883	57.56%	8.093%
53	10.384	1388	1391	1398	rVV	25299	46541	3.25%	0.457%
54	10.445	1398	1401	1403	rVV2	3513	4680	0.33%	0.046%
55	10.500	1404	1410	1417	rVB3	28641	52806	3.68%	0.518%
56	10.579	1417	1423	1430	rBV2	71995	130009	9.07%	1.275%
57	10.664	1432	1437	1439	rVV3	15803	26515	1.85%	0.260%
58	10.701	1439	1443	1448	rVV2	26060	47153	3.29%	0.463%
59	10.762	1448	1453	1462	rVB4	24616	48936	3.41%	0.480%
60	10.859	1462	1469	1474	rBV2	39876	72281	5.04%	0.709%
61	10.920	1474	1479	1492	rVV	162114	292933	20.44%	2.874%
62	11.061	1494	1502	1507	rVV3	33734	65821	4.59%	0.646%
63	11.103	1507	1509	1512	rVV3	10503	16658	1.16%	0.163%
64	11.146	1512	1516	1518	rVV3	20714	35661	2.49%	0.350%
65	11.176	1518	1521	1526	rVV	28444	45703	3.19%	0.448%
66	11.231	1526	1530	1536	rVB4	7306	12768	0.89%	0.125%
67	11.280	1536	1538	1543	rVB6	2795	3638	0.25%	0.036%
68	11.335	1543	1547	1551	rBV5	4627	8855	0.62%	0.087%
69	11.432	1555	1563	1571	rVB3	25942	71012	4.96%	0.697%
70	11.512	1572	1576	1579	rVB2	8989	11648	0.81%	0.114%
71	11.628	1588	1595	1603	rBV	389827	657692	45.90%	6.452%
72	11.731	1605	1612	1621	rVV	217541	366388	25.57%	3.594%
73	11.835	1621	1629	1641	rBV	779307	1433004	100.00%	14.059%
74	12.030	1657	1661	1667	rBV7	3098	6610	0.46%	0.065%
75	12.164	1672	1683	1695	rBV	275946	472472	32.97%	4.635%
76	12.310	1705	1707	1709	rBV3	1344	1507	0.11%	0.015%

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 Title : SFAM01.0

77	12.329	1709	1710	1715	rVV5	3429	4863	0.34%	0.048%
78	12.384	1716	1719	1724	rVV5	2625	5029	0.35%	0.049%
79	12.463	1727	1732	1738	rVB4	8033	14058	0.98%	0.138%
80	12.603	1750	1755	1762	rVB	41782	69693	4.86%	0.684%
81	12.688	1762	1769	1777	rBV	227667	375286	26.19%	3.682%
82	12.865	1795	1798	1804	rBV6	4071	7721	0.54%	0.076%
83	12.932	1806	1809	1816	rVB4	9862	17134	1.20%	0.168%
84	13.005	1819	1821	1822	rBV2	1618	1629	0.11%	0.016%
85	13.018	1822	1823	1824	rBV	1797	963	0.07%	0.009%
86	13.194	1850	1852	1856	rVB5	2330	2029	0.14%	0.020%
87	13.249	1856	1861	1864	rBV4	7938	12987	0.91%	0.127%
88	13.298	1865	1869	1878	rVB7	8562	18869	1.32%	0.185%
89	13.402	1883	1886	1891	rVB6	4653	6436	0.45%	0.063%
90	13.463	1895	1896	1898	rBV2	2108	1317	0.09%	0.013%
91	13.554	1905	1911	1925	rBV	222393	374129	26.11%	3.670%
92	13.755	1940	1944	1947	rBV5	2591	5536	0.39%	0.054%
93	13.853	1953	1960	1966	rBV	203577	343651	23.98%	3.371%
94	14.060	1990	1994	2002	rVB2	23255	42517	2.97%	0.417%
95	14.725	2099	2103	2107	rVB6	4326	6069	0.42%	0.060%
96	15.115	2166	2167	2170	rBV3	1940	1444	0.10%	0.014%
97	15.218	2177	2184	2192	rVB7	10447	24395	1.70%	0.239%
98	15.432	2215	2219	2224	rVB2	16457	25097	1.75%	0.246%
99	15.859	2287	2289	2290	rBV2	2059	1132	0.08%	0.011%
100	15.987	2308	2310	2313	rBV4	1962	2014	0.14%	0.020%

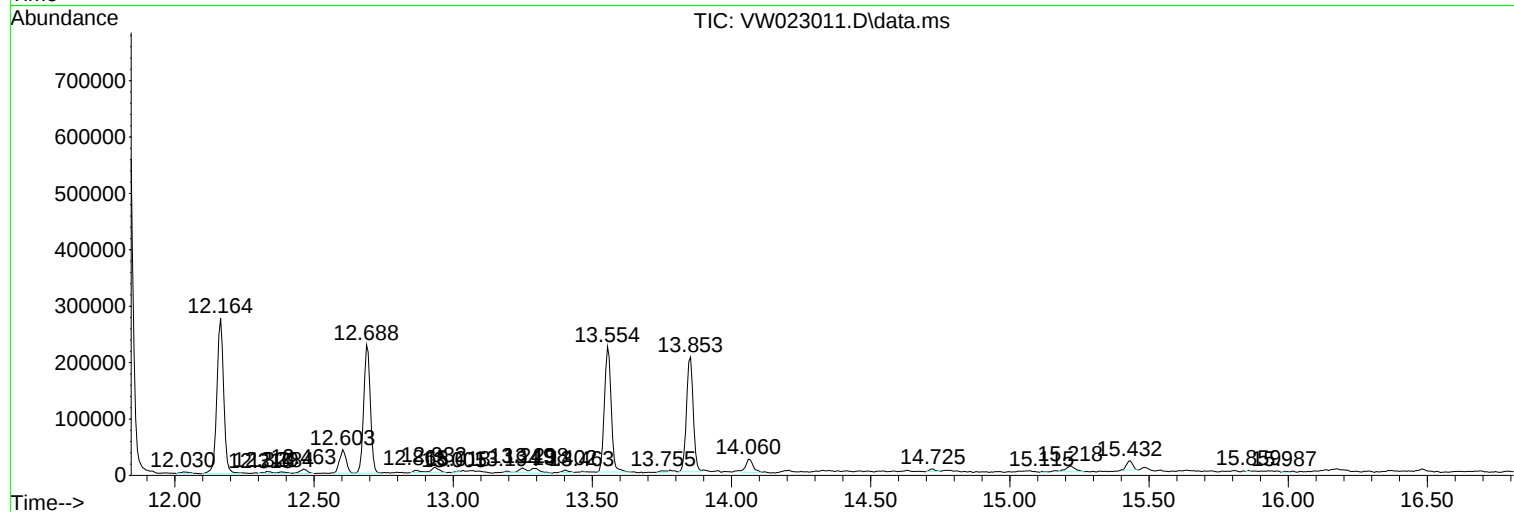
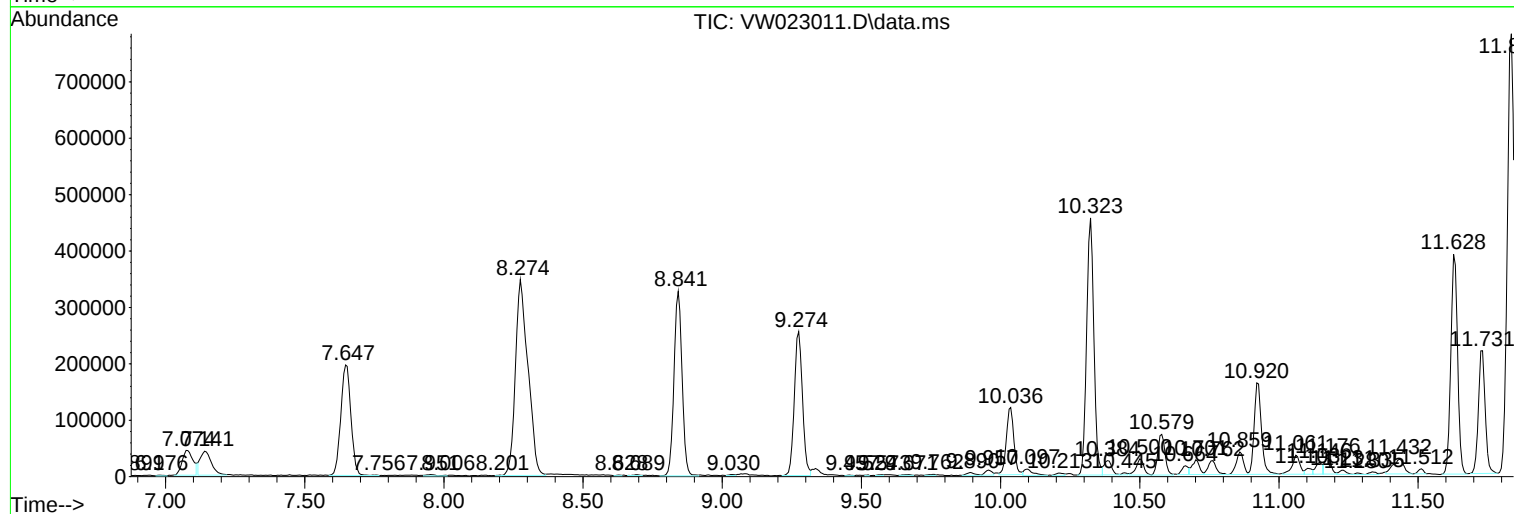
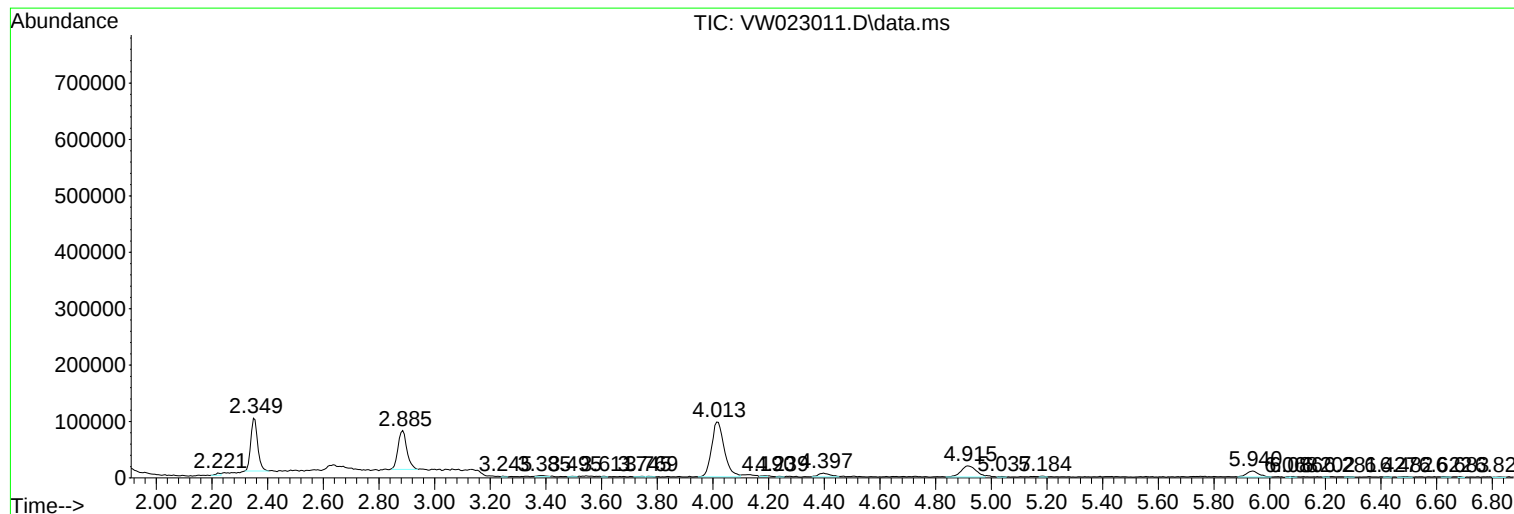
Sum of corrected areas: 10193120

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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM050222SMA.M
Quant Title : SFAM01.0

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



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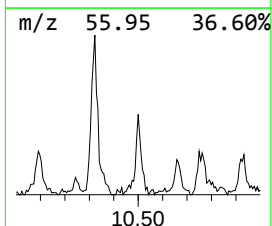
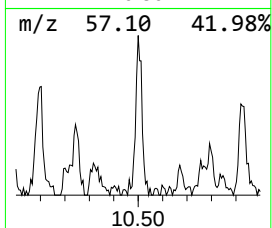
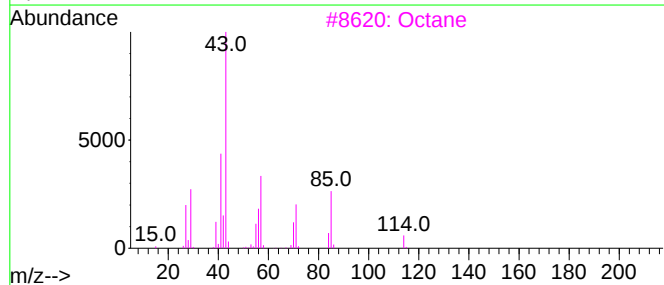
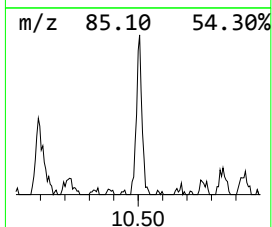
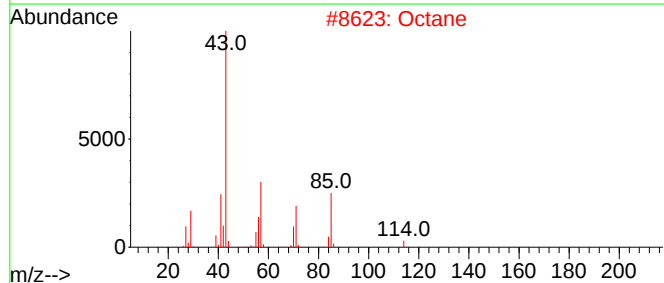
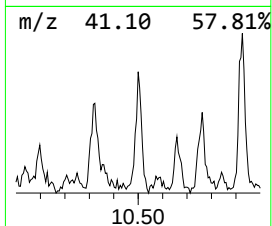
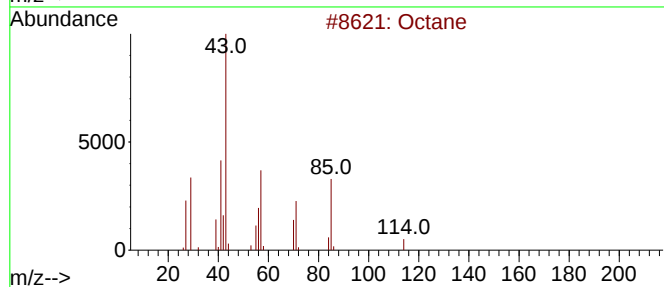
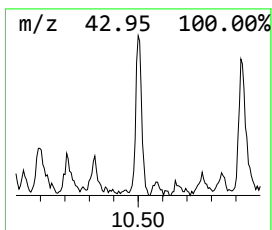
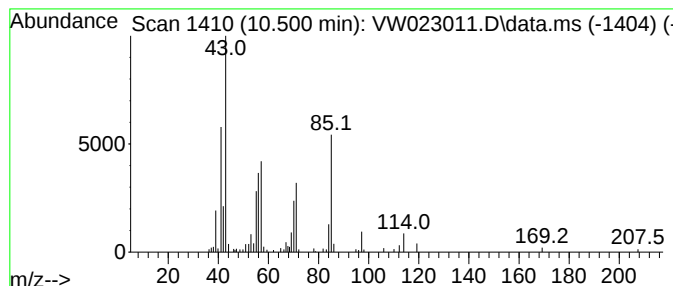
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM050222SMA.M
Quant Title : SFAM01.0

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

Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.500	2.01 ug/L	52806	Chlorobenzene-d5	11.628

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octane			114	C8H18	000111-65-9	87
2	Octane			114	C8H18	000111-65-9	80
3	Octane			114	C8H18	000111-65-9	64
4	Octane			114	C8H18	000111-65-9	59
5	Heptane, 2,4-dimethyl-			128	C9H20	002213-23-2	50



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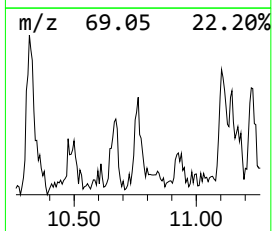
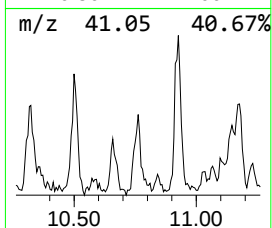
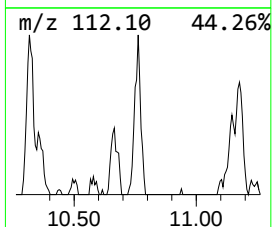
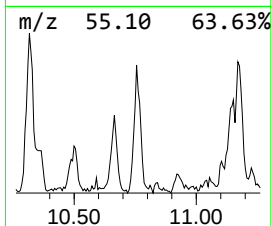
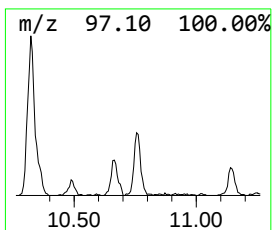
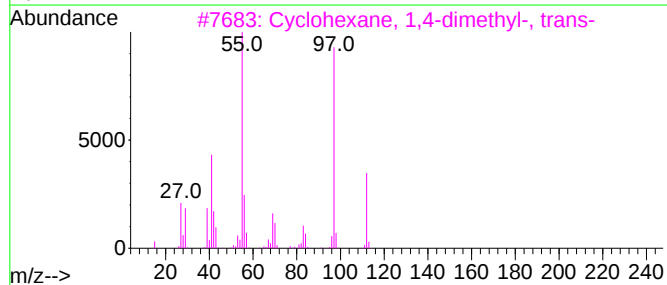
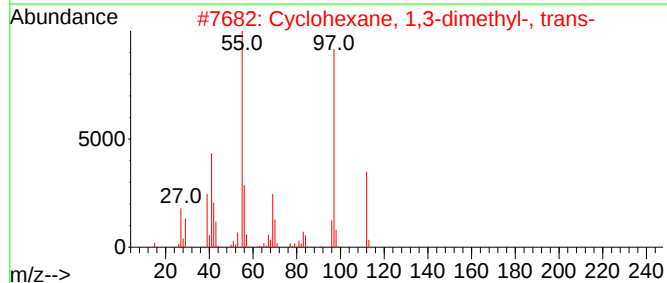
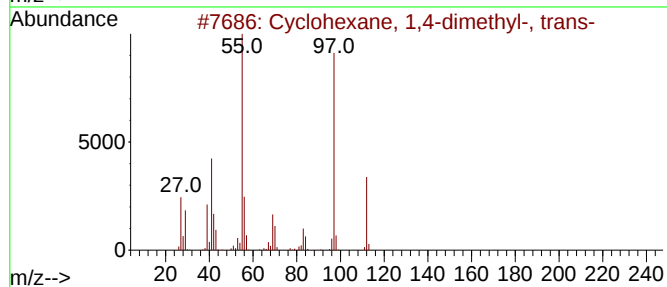
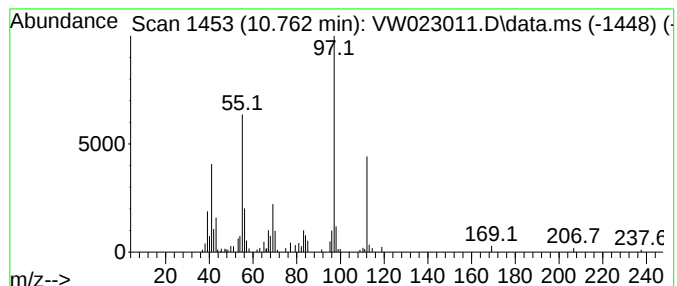
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM050222SMA.M
Quant Title : SFAM01.0

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

Peak Number 5 (DEL) Alkane: Cyclic10.762 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.762	1.86 ug/L	48936	Chlorobenzene-d5	11.628

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, 1,4-dimethyl-, trans-	112	C8H16	002207-04-7	90
2			Cyclohexane, 1,3-dimethyl-, trans-	112	C8H16	002207-03-6	90
3			Cyclohexane, 1,4-dimethyl-, trans-	112	C8H16	002207-04-7	90
4			Cyclohexane, 1,3-dimethyl-, trans-	112	C8H16	002207-03-6	90
5			Cyclohexane, 1,3-dimethyl-, trans-	112	C8H16	002207-03-6	90



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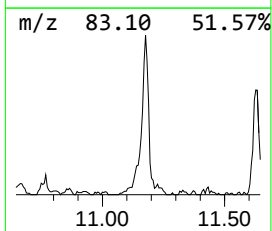
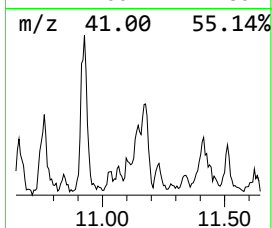
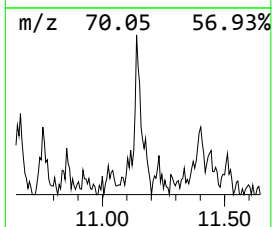
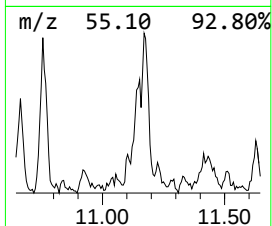
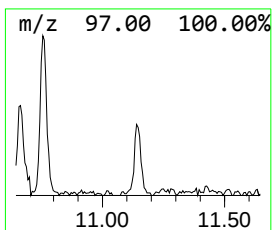
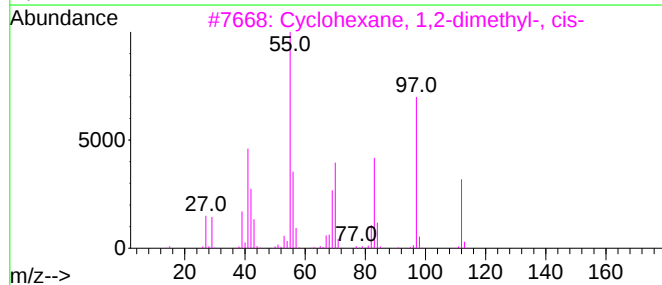
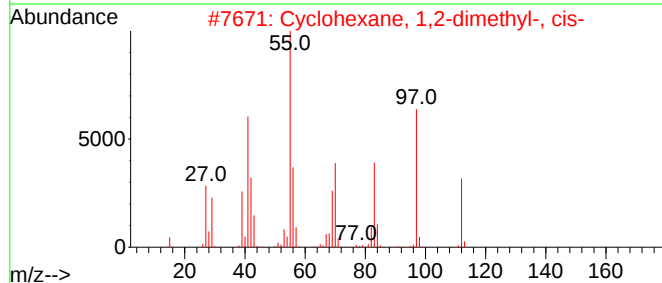
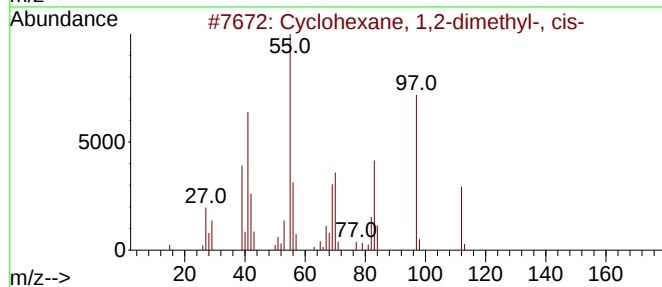
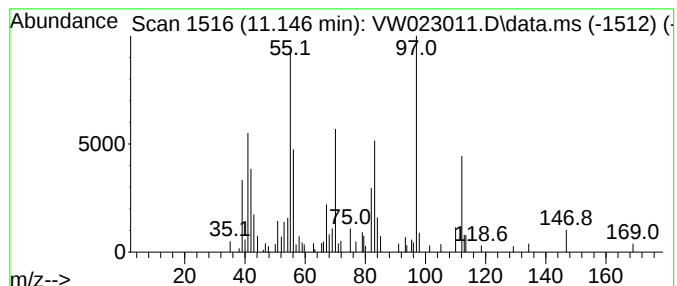
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM050222SMA.M
Quant Title : SFAM01.0

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

Peak Number 7 (DEL) Alkane: Cyclic11.146 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.146	1.36 ug/L	35661	Chlorobenzene-d5	11.628

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, 1,2-dimethyl-, cis-	112	C8H16	002207-01-4	90
2			Cyclohexane, 1,2-dimethyl-, cis-	112	C8H16	002207-01-4	87
3			Cyclohexane, 1,2-dimethyl-, cis-	112	C8H16	002207-01-4	87
4			Cyclohexane, 1,2-dimethyl-, (cis/...	112	C8H16	000583-57-3	87
5			Cyclohexane, 1,2-dimethyl-, (cis/...	112	C8H16	000583-57-3	58



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW051222\
Data File : VW023011.D
Acq On : 12 May 2022 19:06
Operator : SY/VA
Sample : N2767-09
Misc : 4.76g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
BGPM4

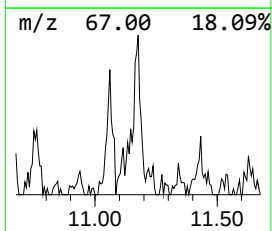
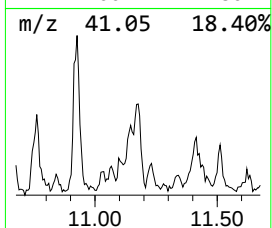
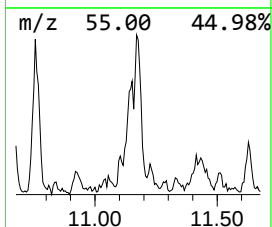
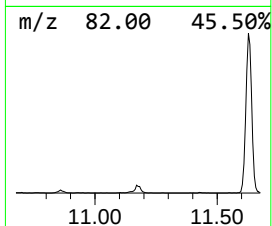
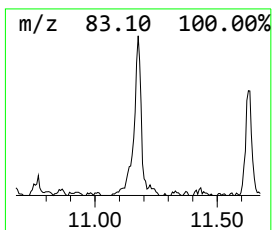
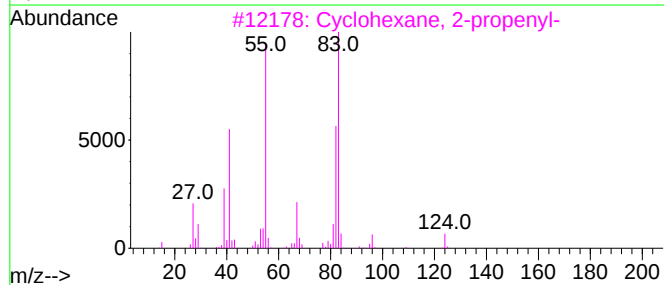
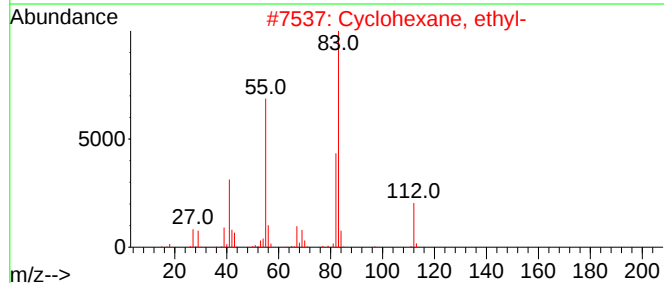
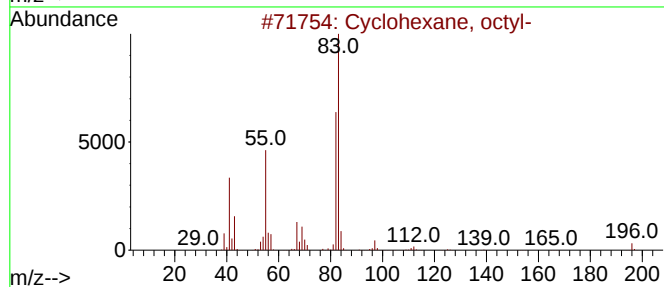
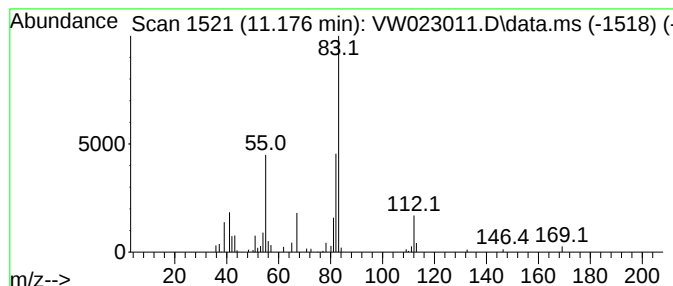
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Quant Title : SFAM01.0

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

Peak Number 8 (DEL) Alkane: Cyclic11.176 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.176	1.74 ug/L	45703	Chlorobenzene-d5	11.628

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, octyl-	196	C14H28	001795-15-9	72
2			Cyclohexane, ethyl-	112	C8H16	001678-91-7	56
3			Cyclohexane, 2-propenyl-	124	C9H16	002114-42-3	56
4			Cyclohexane, ethyl-	112	C8H16	001678-91-7	50
5			Cyclohexane, ethyl-	112	C8H16	001678-91-7	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW051222\
Data File : VW023011.D
Acq On : 12 May 2022 19:06
Operator : SY/VA
Sample : N2767-09
Misc : 4.76g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
BGPM4

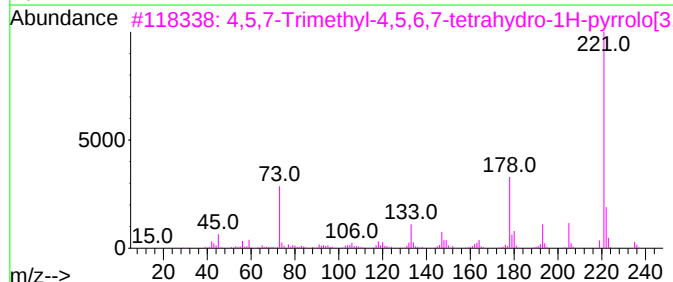
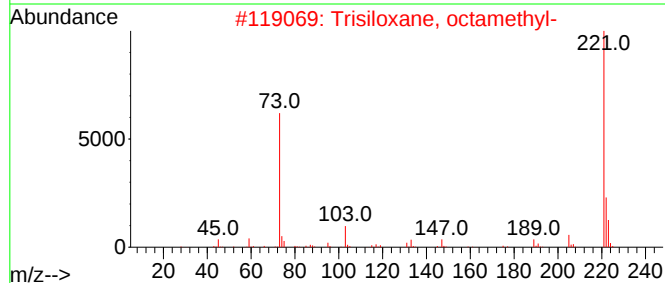
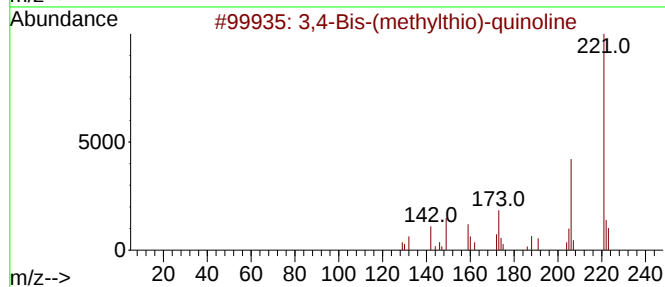
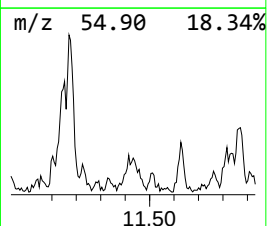
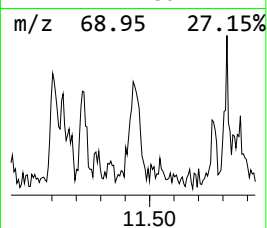
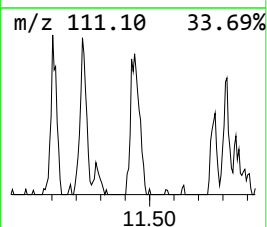
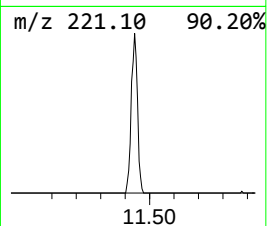
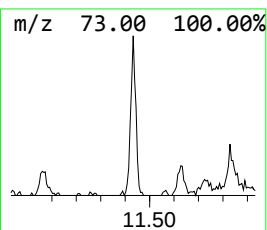
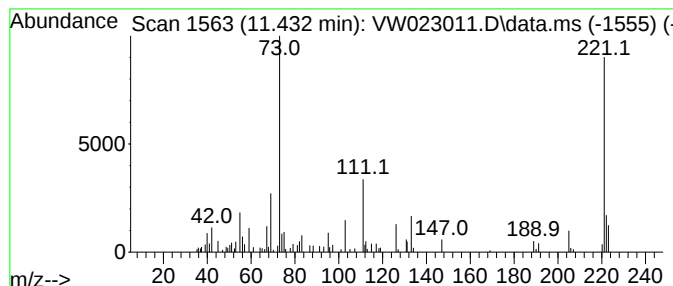
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Quant Title : SFAM01.0

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

Peak Number 9 3,4-Bis-(methylthio)-quinoline Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.432	2.70 ug/L	71012	Chlorobenzene-d5	11.628

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3,4-Bis-(methylthio)-quinoline	221	C11H11NS2	074579-34-3	64
2			Trisiloxane, octamethyl-	236	C8H24O2Si3	000107-51-7	59
3			4,5,7-Trimethyl-4,5,6,7-tetrahyd...	236	C13H24N2Si	1000475-28-6	53
4			1H-Indole, 1-ethyl-2-phenyl-	221	C16H15N	013228-39-2	43
5			4-Isothiazolocarboxylic acid, 3,...	221	C6H7NO2S3	004886-15-1	43



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW051222\
Data File : VW023011.D
Acq On : 12 May 2022 19:06
Operator : SY/VA
Sample : N2767-09
Misc : 4.76g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
BGPM4

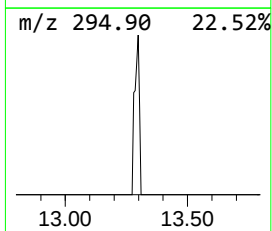
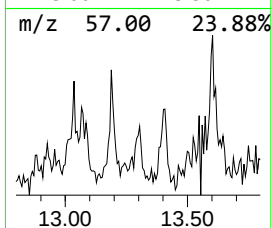
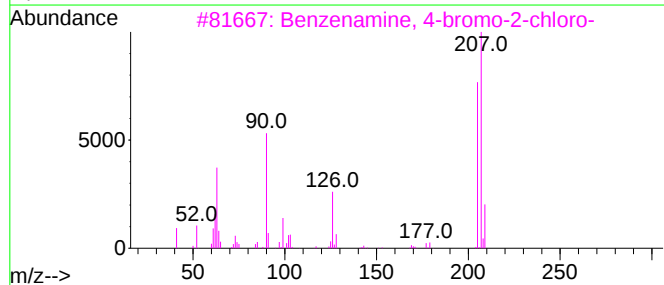
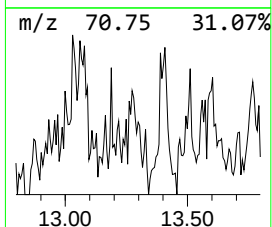
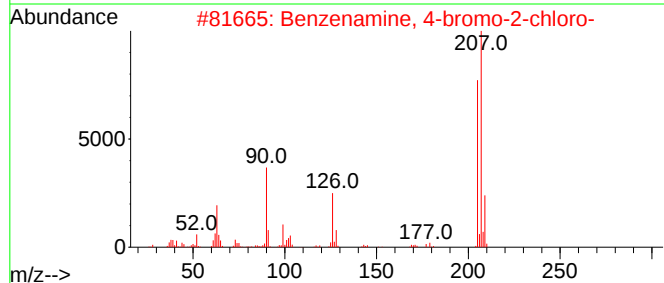
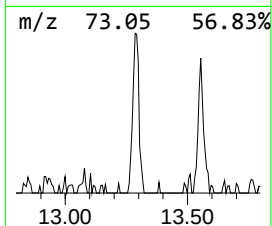
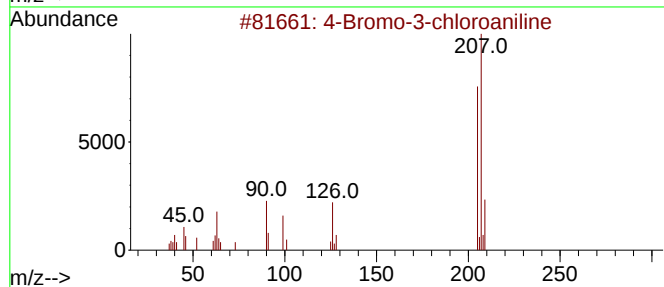
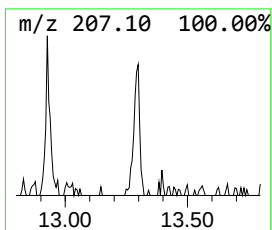
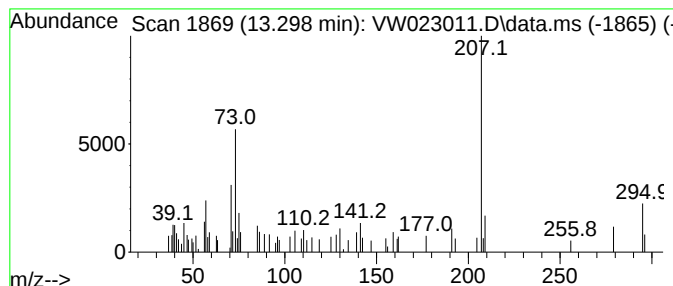
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Quant Title : SFAM01.0

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

Peak Number 11 unknown-01 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.298	1.26 ug/L	18869	1,4-Dichlorobenzene-d4	13.554

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4-Bromo-3-chloroaniline	205	C6H5BrClN	021402-26-6	38
2			Benzenamine, 4-bromo-2-chloro-	205	C6H5BrClN	038762-41-3	38
3			Benzenamine, 4-bromo-2-chloro-	205	C6H5BrClN	038762-41-3	35
4			1,1,1,3,5,5,5-Heptamethyltrisilo...	222	C7H22O2Si3	001873-88-7	32
5			2-Nitro-4-(trifluoromethyl)phenol	207	C7H4F3NO3	000400-99-7	30



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW051222\
Data File : VW023011.D
Acq On : 12 May 2022 19:06
Operator : SY/VA
Sample : N2767-09
Misc : 4.76g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
BGPM4

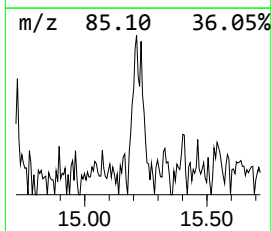
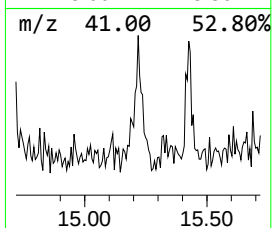
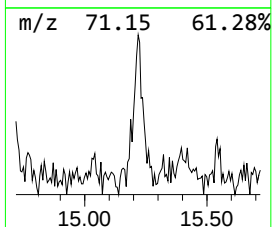
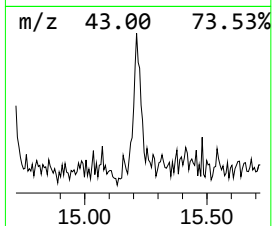
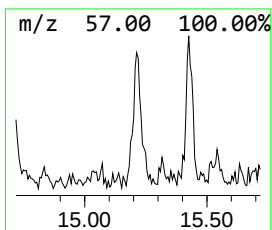
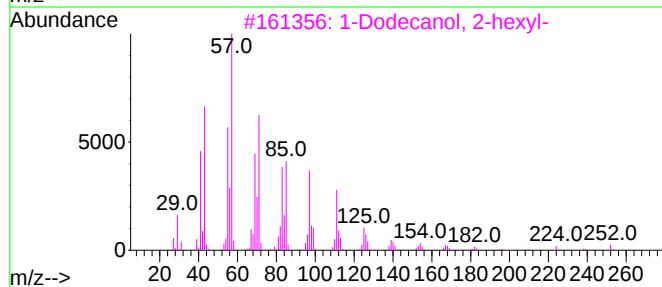
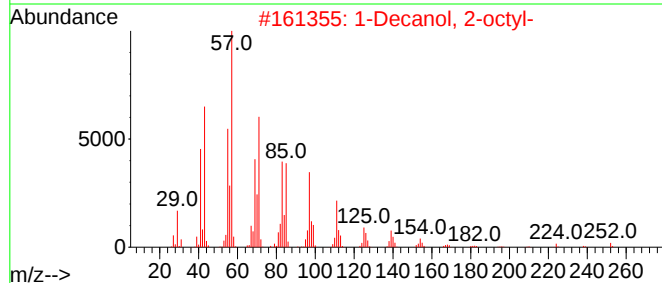
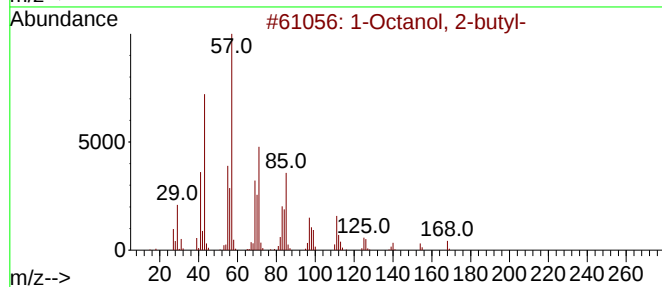
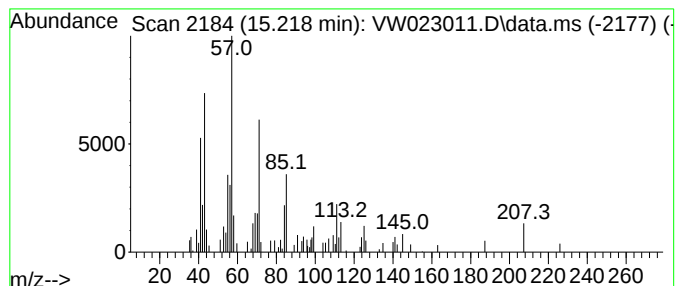
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Quant Title : SFAM01.0

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

Peak Number 13 1-Octanol, 2-butyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.218	1.63 ug/L	24395	1,4-Dichlorobenzene-d4	13.554

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1-Octanol, 2-butyl-	186	C12H26O	003913-02-8	80
2		1-Decanol, 2-octyl-	270	C18H38O	045235-48-1	72
3		1-Dodecanol, 2-hexyl-	270	C18H38O	110225-00-8	72
4		1-Decanol, 2-hexyl-	242	C16H34O	002425-77-6	72
5		2-Hexyl-1-octanol	214	C14H30O	019780-79-1	72



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW051222\
Data File : VW023011.D
Acq On : 12 May 2022 19:06
Operator : SY/VA
Sample : N2767-09
Misc : 4.76g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
BGPM4

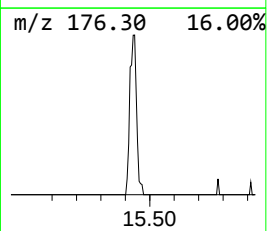
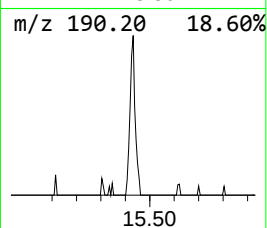
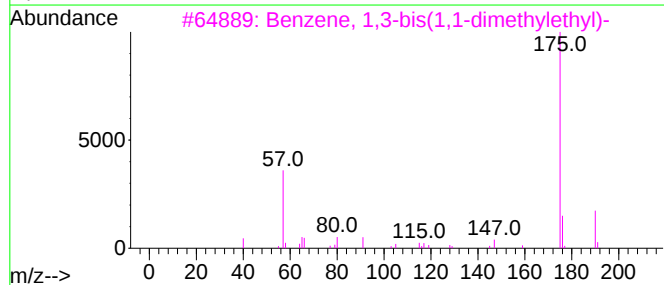
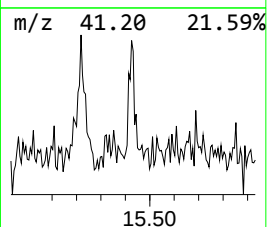
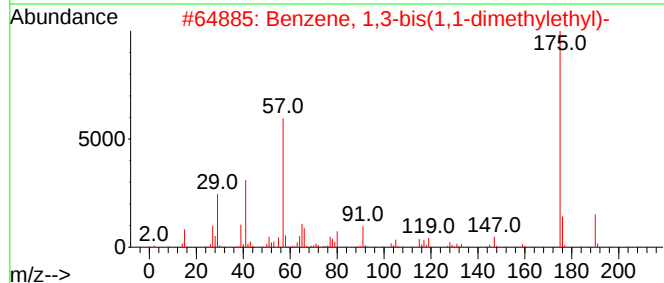
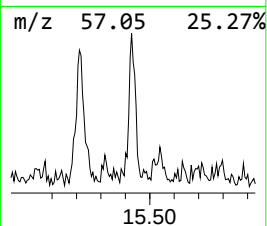
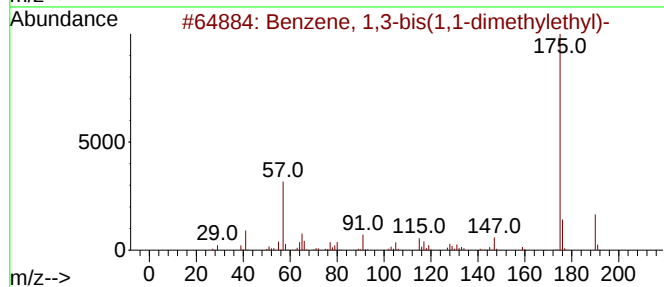
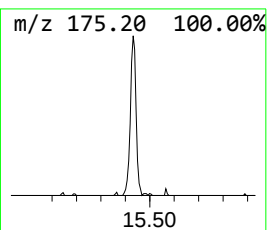
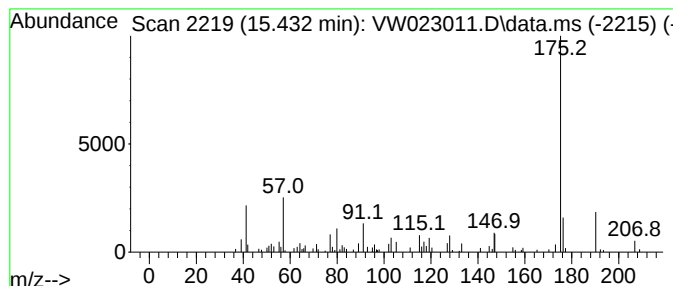
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Quant Title : SFAM01.0

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

Peak Number 14 Benzene, 1,3-bis(1,1-dimeth... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.432	1.68 ug/L	25097	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	72
2			Benzene, 1,3-bis(1,1-dimethyleth...	190	C14H22	001014-60-4	72
3			Benzene, 1,3-bis(1,1-dimethyleth...	190	C14H22	001014-60-4	72
4			Benzene, 1,3-bis(1,1-dimethyleth...	190	C14H22	001014-60-4	68
5			Benzenepropanal, 4-(1,1-dimethyl...	190	C13H18O	018127-01-0	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW051222\
Data File : VW023011.D
Acq On : 12 May 2022 19:06
Operator : SY/VA
Sample : N2767-09
Misc : 4.76g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
BGPM4

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM050222SMA.M
Quant Title : SFAM01.0

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

					--Internal Standard--			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
(DEL) Alkane: S...	10.500	2.0	ug/L	52806	2	11.628	657692	25.0
(DEL) Alkane: C...	10.762	1.9	ug/L	48936	2	11.628	657692	25.0
(DEL) Alkane: C...	11.146	1.4	ug/L	35661	2	11.628	657692	25.0
(DEL) Alkane: C...	11.176	1.7	ug/L	45703	2	11.628	657692	25.0
3,4-Bis-(methyl...	11.432	2.7	ug/L	71012	2	11.628	657692	25.0
unknown-01	13.298	1.3	ug/L	18869	3	13.554	374129	25.0
1-Octanol, 2-bu...	15.218	1.6	ug/L	24395	3	13.554	374129	25.0
Benzene, 1,3-bi...	15.432	1.7	ug/L	25097	3	13.554	374129	25.0