

Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW041222\
 Data File : VW022503.D
 Acq On : 12 Apr 2022 13:33
 Operator : SY/VA
 Sample : N2316-12RE
 Misc : 7.42g/10mL/MSVOA_W/SOIL
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 ETNR2RE

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: VW022503.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.959	5	9	17	rVB	126668	176300	83.57%	8.270%
2	2.885	157	161	166	rVB	13563	23843	11.30%	1.118%
3	3.135	200	202	203	rBV2	938	677	0.32%	0.032%
4	3.398	242	245	250	rBV5	1824	3379	1.60%	0.158%
5	3.495	258	261	263	rBV4	1412	1215	0.58%	0.057%
6	3.891	324	326	329	rVB3	1107	1066	0.51%	0.050%
7	3.934	329	333	337	rBV5	1740	3055	1.45%	0.143%
8	3.971	337	339	340	rVV2	1868	1546	0.73%	0.073%
9	4.026	340	348	358	rVB2	25726	68116	32.29%	3.195%
10	4.379	403	406	407	rBV2	1846	1981	0.94%	0.093%
11	4.446	412	417	419	rVV5	1019	1957	0.93%	0.092%
12	4.495	423	425	427	rVB2	783	586	0.28%	0.027%
13	4.599	439	442	446	rBV4	767	1482	0.70%	0.070%
14	4.678	453	455	457	rVB3	911	585	0.28%	0.027%
15	4.696	457	458	460	rVB2	1076	564	0.27%	0.026%
16	4.721	460	462	464	rBV2	1329	1644	0.78%	0.077%
17	4.751	466	467	471	rVV3	881	620	0.29%	0.029%
18	4.849	482	483	486	rBV3	731	756	0.36%	0.035%
19	5.141	530	531	534	rVB2	995	619	0.29%	0.029%
20	5.202	539	541	545	rVB4	1190	971	0.46%	0.046%
21	5.233	545	546	548	rBV2	789	576	0.27%	0.027%
22	5.416	574	576	577	rBV2	833	702	0.33%	0.033%
23	5.513	590	592	594	rBV3	737	917	0.43%	0.043%
24	5.574	594	602	608	rVB8	3581	11265	5.34%	0.528%
25	5.696	620	622	623	rBV	733	539	0.26%	0.025%
26	5.775	634	635	637	rBV3	767	603	0.29%	0.028%
27	5.873	647	651	652	rBV2	949	1188	0.56%	0.056%
28	5.891	652	654	655	rBV2	810	533	0.25%	0.025%
29	5.909	655	657	658	rBV	750	559	0.26%	0.026%
30	5.940	660	662	664	rBV2	1312	1132	0.54%	0.053%
31	6.324	723	725	729	rBV5	963	1221	0.58%	0.057%
32	6.397	735	737	740	rVB3	831	670	0.32%	0.031%
33	6.495	752	753	755	rBV2	941	545	0.26%	0.026%
34	6.519	755	757	759	rBV2	712	753	0.36%	0.035%
35	6.629	772	775	778	rBV5	1214	1794	0.85%	0.084%
36	6.702	785	787	790	rVV4	1025	1099	0.52%	0.052%

Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW041222\
 Data File : VW022503.D
 Acq On : 12 Apr 2022 13:33
 Operator : SY/VA
 Sample : N2316-12RE
 Misc : 7.42g/10mL/MSVOA_W/SOIL
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 ETNR2RE

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

37	6.793	800	802	803	rVB2	832	537	0.25%	0.025%
38	6.812	803	805	808	rBV4	690	980	0.46%	0.046%
39	6.866	812	814	816	rBV3	942	854	0.40%	0.040%
40	6.970	825	831	833	rBV6	2832	4552	2.16%	0.214%
41	7.092	843	851	855	rBV3	6547	18936	8.98%	0.888%
42	7.251	874	877	878	rBV3	1555	1406	0.67%	0.066%
43	7.287	881	883	884	rBV2	1165	1165	0.55%	0.055%
44	7.330	888	890	892	rBV3	991	1161	0.55%	0.054%
45	7.391	898	900	904	rVB4	1245	1072	0.51%	0.050%
46	7.433	904	907	910	rBV4	788	1045	0.50%	0.049%
47	7.464	910	912	914	rVB3	1052	781	0.37%	0.037%
48	7.494	914	917	918	rBV	1068	1017	0.48%	0.048%
49	7.580	929	931	933	rBV3	740	832	0.39%	0.039%
50	7.653	935	943	954	rVB2	34008	89027	42.20%	4.176%
51	7.909	983	985	987	rBV3	911	817	0.39%	0.038%
52	7.958	988	993	999	rVB7	3281	6704	3.18%	0.314%
53	8.147	1023	1024	1027	rBV3	723	783	0.37%	0.037%
54	8.171	1027	1028	1032	rVB4	1405	1236	0.59%	0.058%
55	8.214	1032	1035	1036	rBV3	983	875	0.41%	0.041%
56	8.281	1036	1046	1047	rBV	63732	119989	56.88%	5.628%
57	8.695	1112	1114	1116	rVV3	1663	1407	0.67%	0.066%
58	8.714	1116	1117	1119	rVV2	1648	1016	0.48%	0.048%
59	8.738	1119	1121	1123	rVB3	848	632	0.30%	0.030%
60	8.842	1131	1138	1147	rBV	60454	123875	58.72%	5.811%
61	9.006	1164	1165	1166	rBV	1136	506	0.24%	0.024%
62	9.092	1175	1179	1185	rVB3	13214	25201	11.95%	1.182%
63	9.275	1201	1209	1216	rBV2	45104	97016	45.99%	4.551%
64	9.457	1237	1239	1242	rBV4	753	1086	0.51%	0.051%
65	9.512	1242	1248	1256	rVB5	6689	15480	7.34%	0.726%
66	9.604	1261	1263	1264	rBV2	773	715	0.34%	0.034%
67	9.738	1283	1285	1289	rVB4	1243	1354	0.64%	0.064%
68	9.781	1289	1292	1294	rBV3	703	889	0.42%	0.042%
69	9.805	1294	1296	1298	rBV3	973	634	0.30%	0.030%
70	9.957	1315	1321	1327	rVV5	9741	19583	9.28%	0.919%
71	10.037	1328	1334	1340	rVB	22752	41684	19.76%	1.955%
72	10.323	1373	1381	1387	rBV	89481	161540	76.57%	7.577%
73	10.384	1387	1391	1398	rVV2	22636	40963	19.42%	1.921%
74	10.457	1401	1403	1406	rVB4	1506	1133	0.54%	0.053%
75	10.482	1406	1407	1408	rBV	1503	951	0.45%	0.045%
76	10.579	1418	1423	1429	rVB2	11989	23206	11.00%	1.089%

Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW041222\
 Data File : VW022503.D
 Acq On : 12 Apr 2022 13:33
 Operator : SY/VA
 Sample : N2316-12RE
 Misc : 7.42g/10mL/MSVOA_W/SOIL
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 ETNR2RE

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

77	10.713	1441	1445	1450	rBV6	4338	8690	4.12%	0.408%
78	10.927	1474	1480	1492	rBV	112957	210965	100.00%	9.896%
79	11.280	1536	1538	1539	rBV2	1349	1378	0.65%	0.065%
80	11.366	1550	1552	1554	rVB2	1305	935	0.44%	0.044%
81	11.390	1554	1556	1557	rBV2	929	740	0.35%	0.035%
82	11.463	1566	1568	1569	rBV2	1131	770	0.36%	0.036%
83	11.634	1588	1596	1598	rBV	89553	176434	83.63%	8.276%
84	12.164	1680	1683	1687	rVB6	3511	5664	2.68%	0.266%
85	12.256	1696	1698	1699	rVB2	911	509	0.24%	0.024%
86	12.286	1699	1703	1704	rBV4	1258	1613	0.76%	0.076%
87	12.689	1763	1769	1773	rBV2	37511	65438	31.02%	3.070%
88	12.743	1773	1778	1784	rVB	95968	158467	75.12%	7.433%
89	12.890	1797	1802	1812	rVB2	17363	33561	15.91%	1.574%
90	12.987	1815	1818	1825	rVB2	10853	15675	7.43%	0.735%
91	13.103	1832	1837	1843	rVB2	12399	22424	10.63%	1.052%
92	13.170	1846	1848	1849	rBV2	1088	703	0.33%	0.033%
93	13.554	1905	1911	1922	rVB	82484	144746	68.61%	6.790%
94	13.847	1953	1959	1969	rVB	75628	137395	65.13%	6.445%
95	14.091	1997	1999	2005	rVB5	3860	5099	2.42%	0.239%
96	14.469	2057	2061	2066	rVB7	5352	8672	4.11%	0.407%
97	14.749	2105	2107	2108	rBV2	1184	754	0.36%	0.035%
98	14.993	2143	2147	2149	rVB4	1239	1417	0.67%	0.066%
99	15.926	2298	2300	2302	rVB3	1627	913	0.43%	0.043%
100	16.523	2396	2398	2400	rBV3	1128	1206	0.57%	0.057%

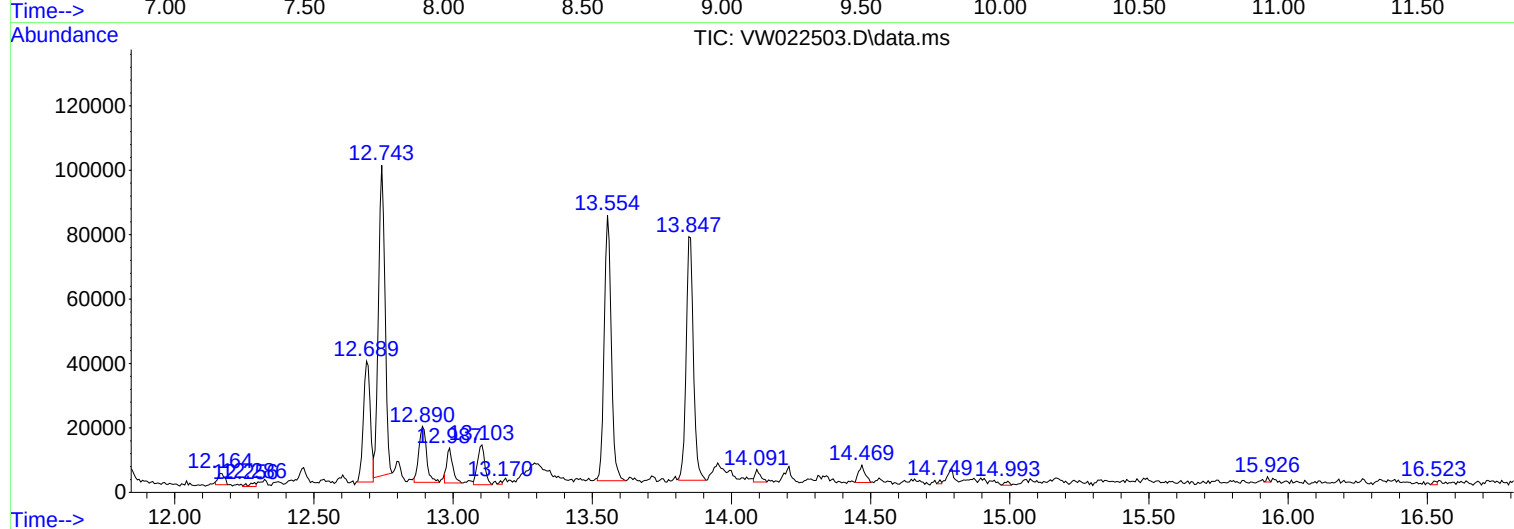
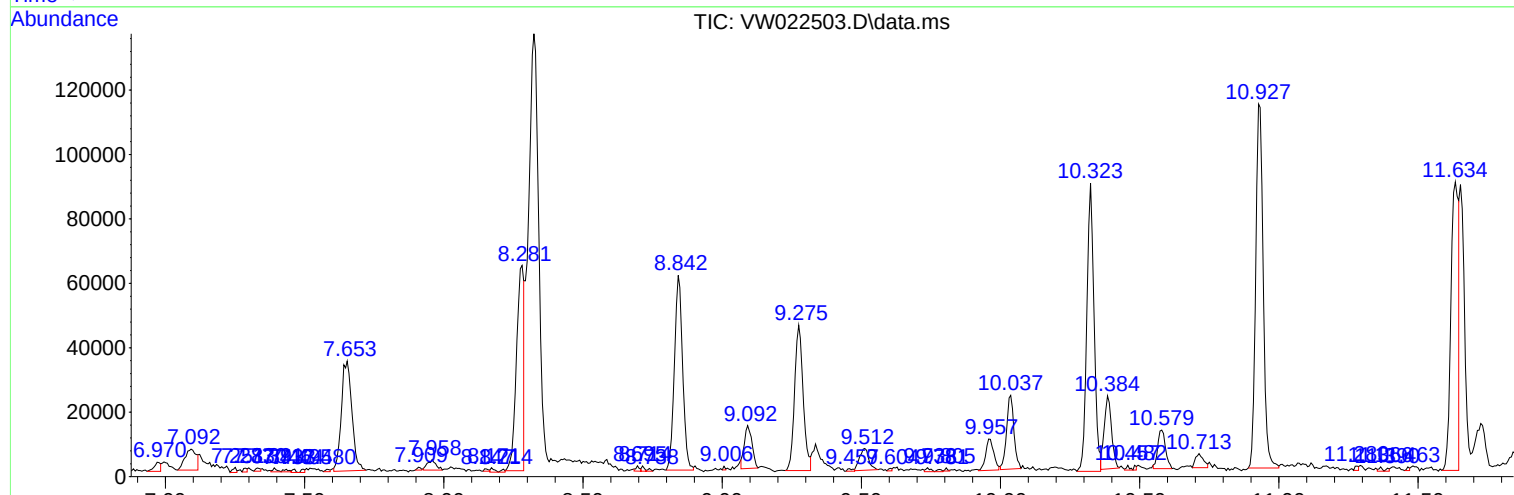
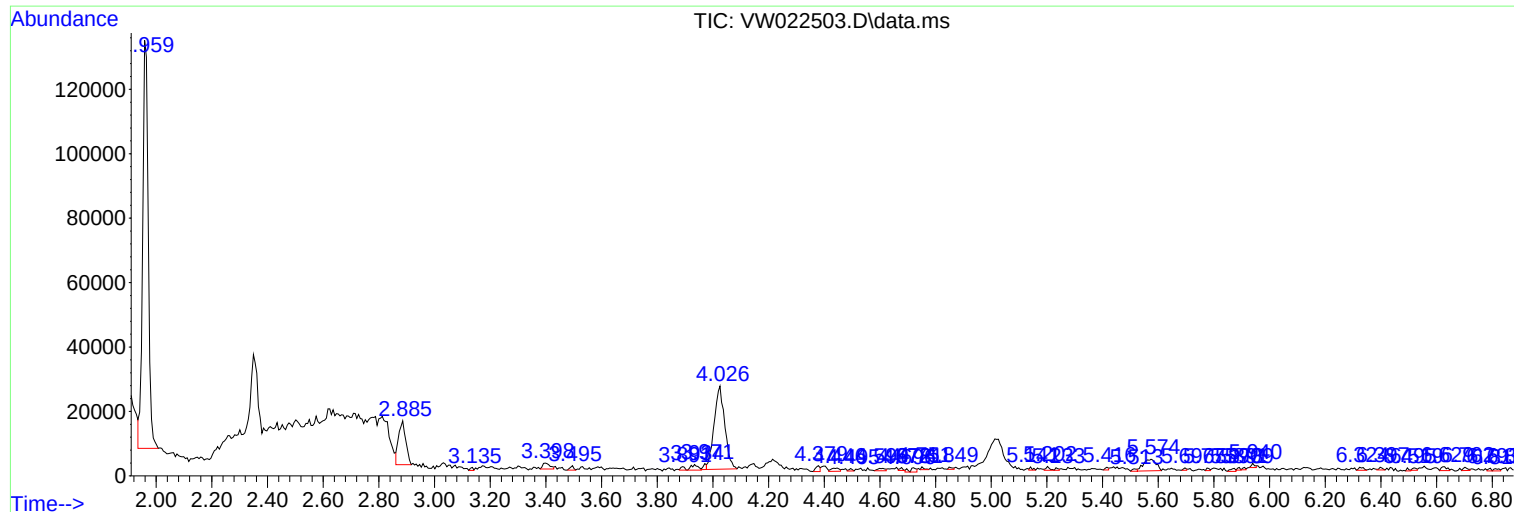
Sum of corrected areas: 2131866

Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW041222\
Data File : VW022503.D
Acq On : 12 Apr 2022 13:33
Operator : SY/VA
Sample : N2316-12RE
Misc : 7.42g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
ETNR2RE

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW041222\
Data File : VW022503.D
Acq On : 12 Apr 2022 13:33
Operator : SY/VA
Sample : N2316-12RE
Misc : 7.42g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
ETNR2RE

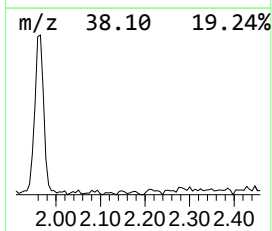
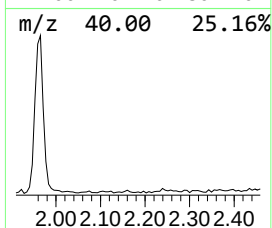
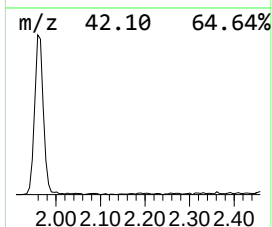
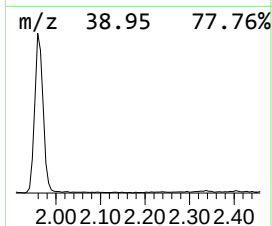
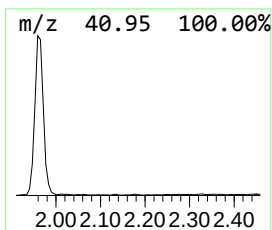
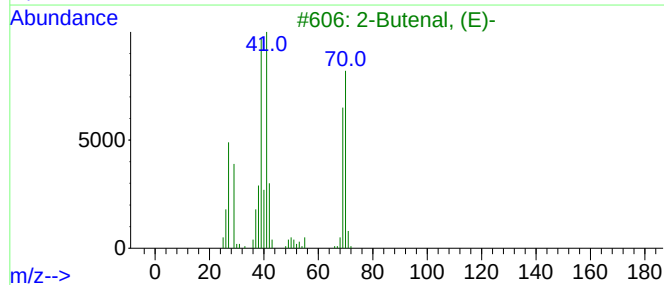
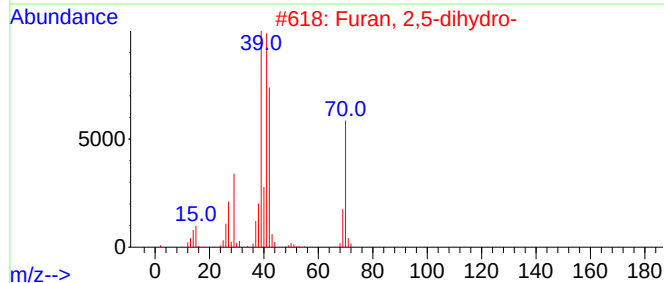
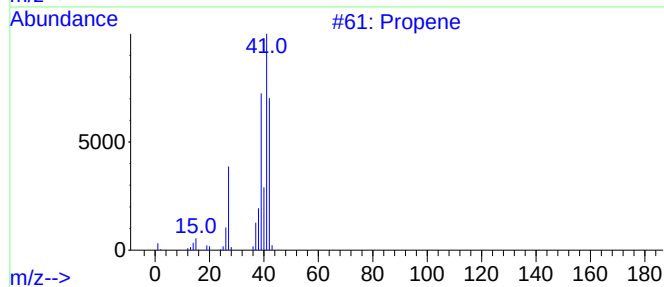
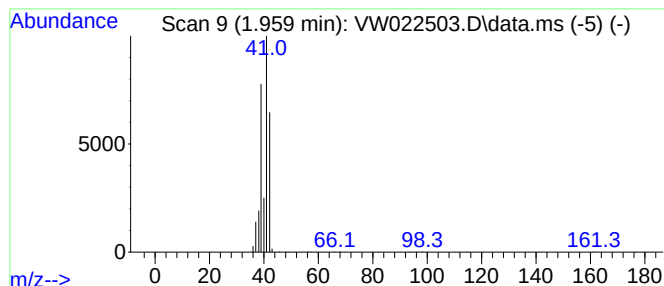
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM040722SMA.M
Quant Title : SFAM01.0

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

Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.959	35.58 ug/L	176300	1,4-Difluorobenzene	8.842

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propene	42	C3H6	000115-07-1	90
2			Furan, 2,5-dihydro-	70	C4H6O	001708-29-8	83
3			2-Butenal, (E)-	70	C4H6O	000123-73-9	78
4			3-Butenoic acid	86	C4H6O2	000625-38-7	78
5			N-Methyl-7-azabicyclo(2,2,1)hept...	109	C7H11N	055590-26-6	74



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW041222\
Data File : VW022503.D
Acq On : 12 Apr 2022 13:33
Operator : SY/VA
Sample : N2316-12RE
Misc : 7.42g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
ETNR2RE

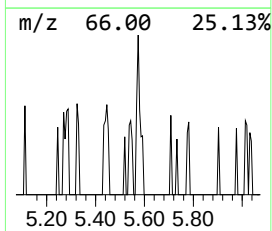
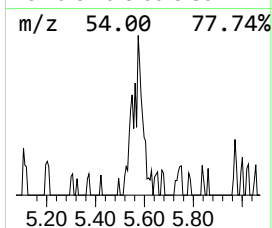
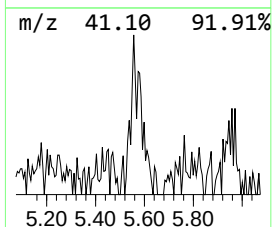
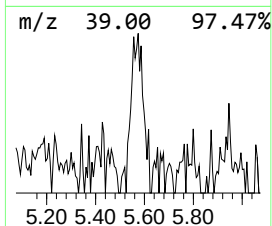
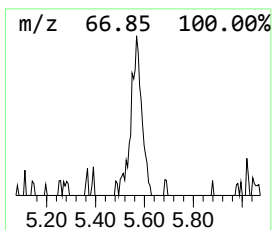
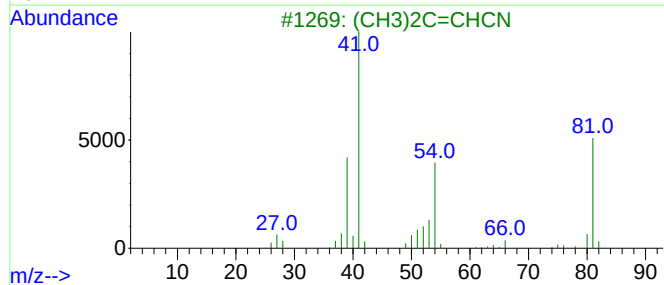
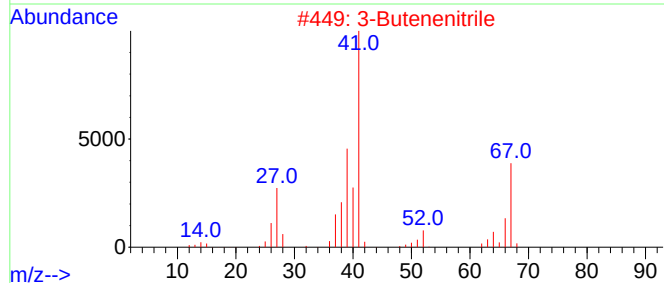
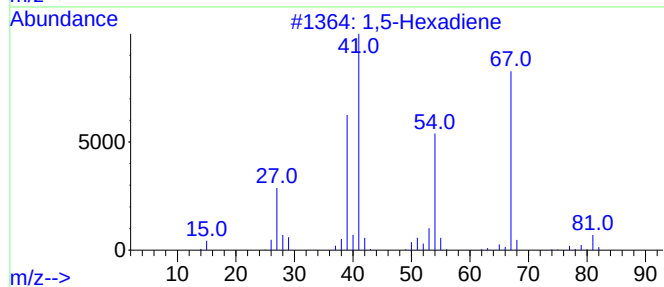
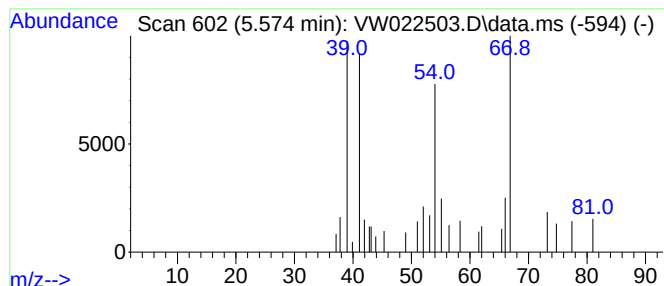
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM040722SMA.M
Quant Title : SFAM01.0

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

Peak Number 2 unknown-01 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.574	2.27 ug/L	11265	1,4-Difluorobenzene	8.842

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1,5-Hexadiene	82	C6H10	000592-42-7	45
2		3-Butenenitrile	67	C4H5N	000109-75-1	43
3		(CH3)2C=CHCN	81	C5H7N	004786-24-7	38
4		2-Butenenitrile	67	C4H5N	004786-20-3	37
5		Bicyclo[1.1.0]butane	54	C4H6	000157-33-5	33



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW041222\
Data File : VW022503.D
Acq On : 12 Apr 2022 13:33
Operator : SY/VA
Sample : N2316-12RE
Misc : 7.42g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
ETNR2RE

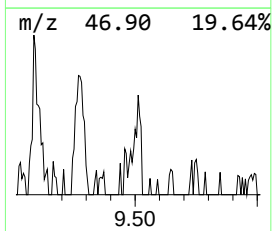
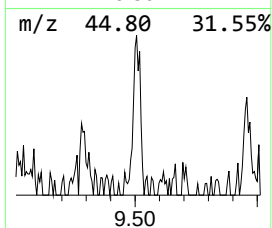
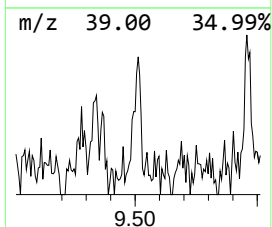
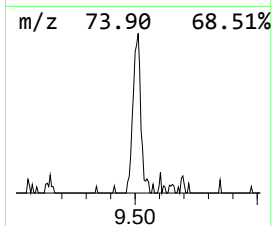
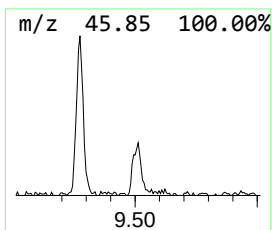
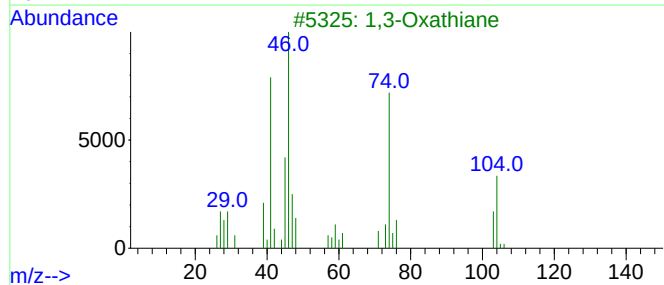
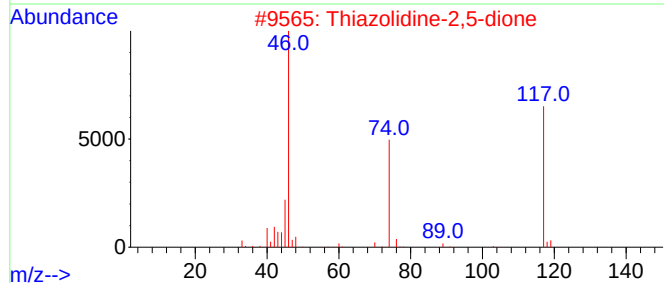
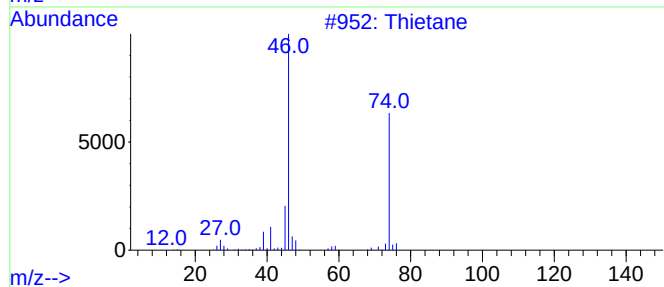
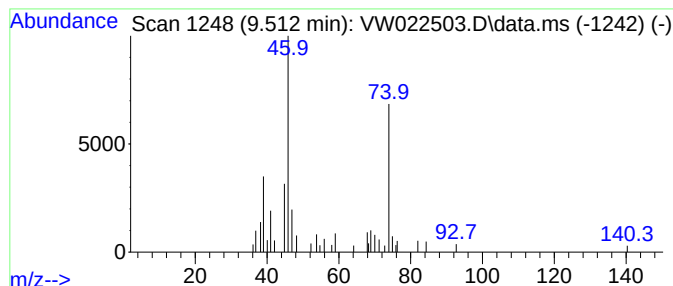
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM040722SMA.M
Quant Title : SFAM01.0

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

Peak Number 3 Thietane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.512	3.12 ug/L	15480	1,4-Difluorobenzene	8.842

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Thietane	74	C3H6S	000287-27-4	50
2			Thiazolidine-2,5-dione	117	C3H3N02S	016874-97-8	40
3			1,3-Oxathiane	104	C4H8OS	000646-12-8	25
4			Thietane, 3-methyl-	88	C4H8S	022438-40-0	23
5			Thiazolidin-4-one, 2-hydroxymeth...	146	C4H6N2O2S	104959-22-0	23



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW041222\
Data File : VW022503.D
Acq On : 12 Apr 2022 13:33
Operator : SY/VA
Sample : N2316-12RE
Misc : 7.42g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
ETNR2RE

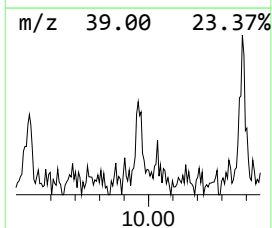
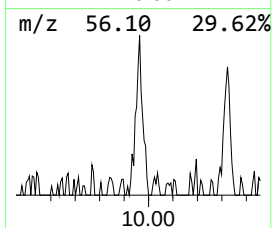
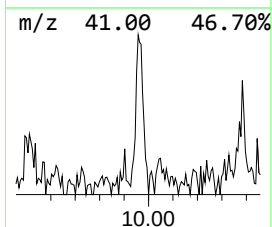
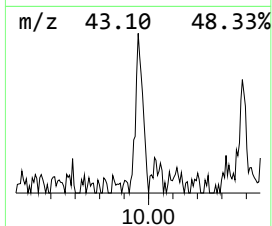
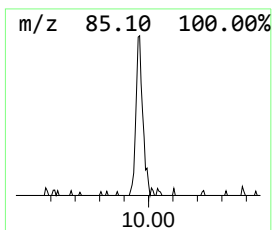
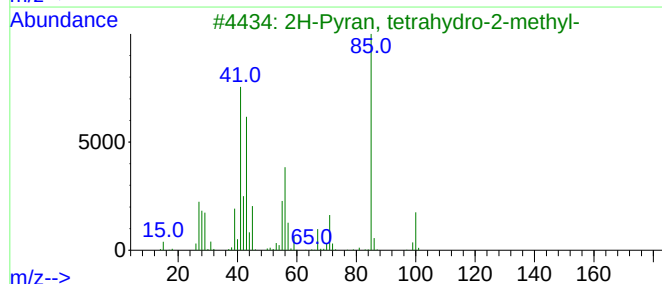
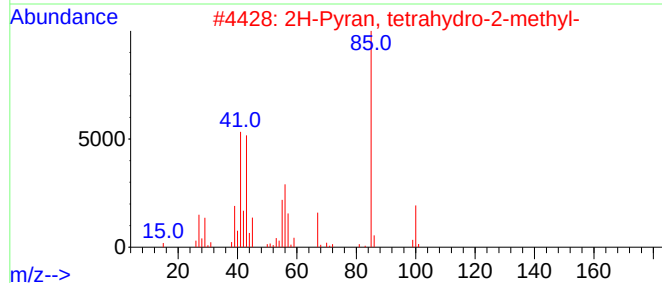
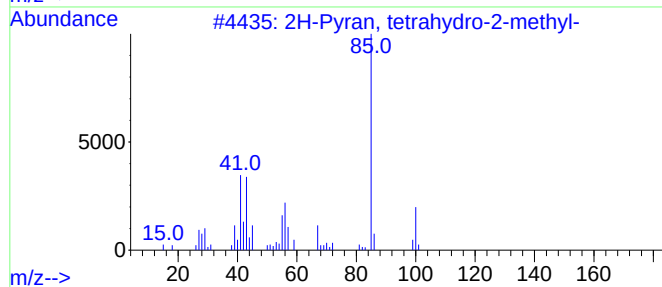
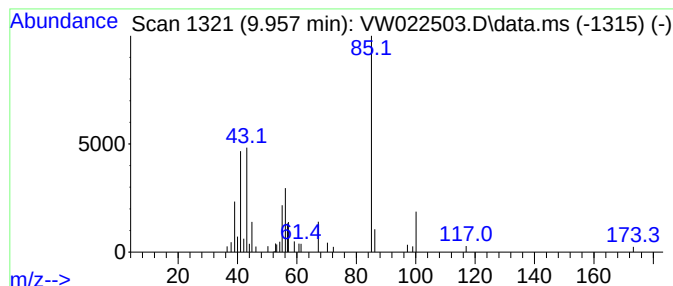
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM040722SMA.M
Quant Title : SFAM01.0

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

Peak Number 4 2H-Pyran, tetrahydro-2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.957	3.95 ug/L	19583	1,4-Difluorobenzene	8.842

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2H-Pyran, tetrahydro-2-methyl-			100	C6H12O	010141-72-7	86
2	2H-Pyran, tetrahydro-2-methyl-			100	C6H12O	010141-72-7	83
3	2H-Pyran, tetrahydro-2-methyl-			100	C6H12O	010141-72-7	72
4	1-Bromo-8-tetrahydropyranyloxyoc...			292	C13H25BrO2	050816-20-1	59
5	Furfuryl alcohol, tetrahydro-5-m...			116	C6H12O2	016015-08-0	59



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW041222\
Data File : VW022503.D
Acq On : 12 Apr 2022 13:33
Operator : SY/VA
Sample : N2316-12RE
Misc : 7.42g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
ETNR2RE

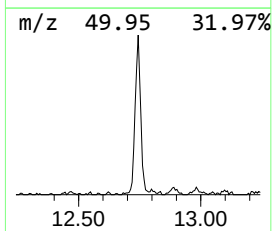
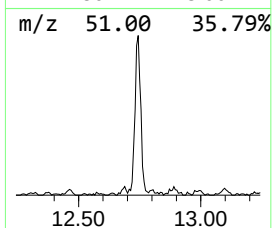
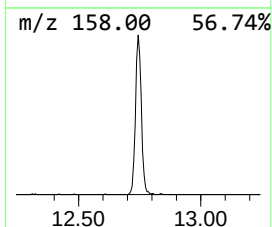
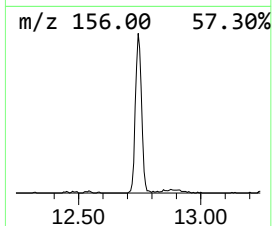
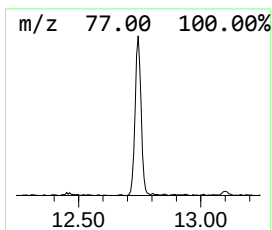
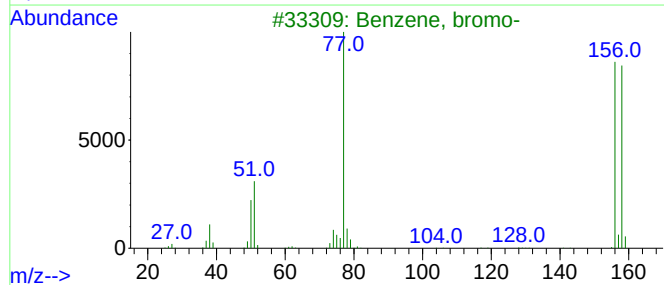
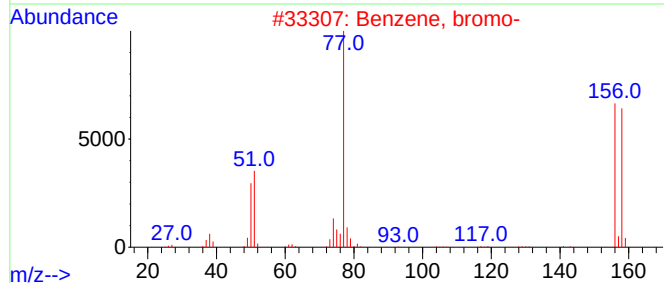
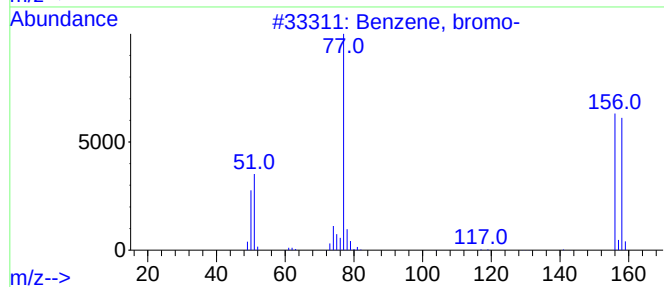
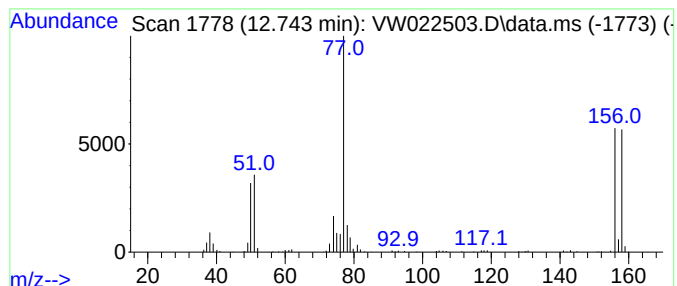
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM040722SMA.M
Quant Title : SFAM01.0

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

Peak Number 6 Benzene, bromo- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.743	27.37 ug/L	158467	1,4-Dichlorobenzene-d4	13.554

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, bromo-	156	C6H5Br	000108-86-1	95
2			Benzene, bromo-	156	C6H5Br	000108-86-1	94
3			Benzene, bromo-	156	C6H5Br	000108-86-1	94
4			Benzene, bromo-	156	C6H5Br	000108-86-1	91
5			Benzene, bromo-	156	C6H5Br	000108-86-1	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW041222\
Data File : VW022503.D
Acq On : 12 Apr 2022 13:33
Operator : SY/VA
Sample : N2316-12RE
Misc : 7.42g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
ETNR2RE

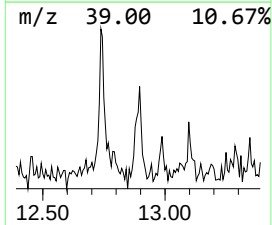
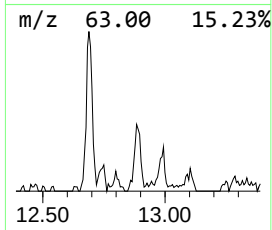
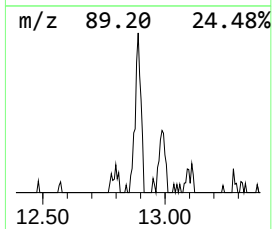
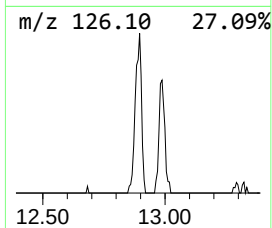
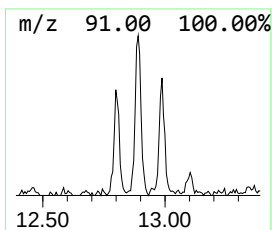
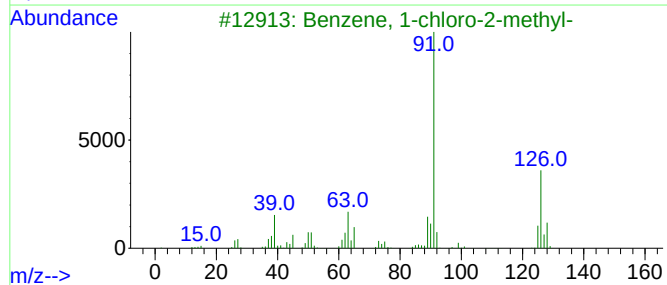
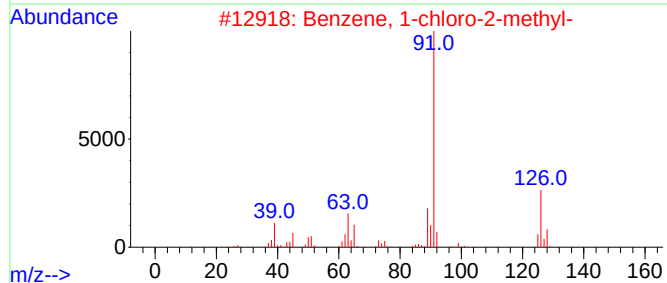
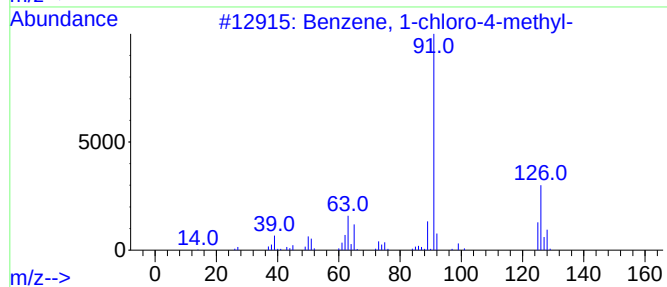
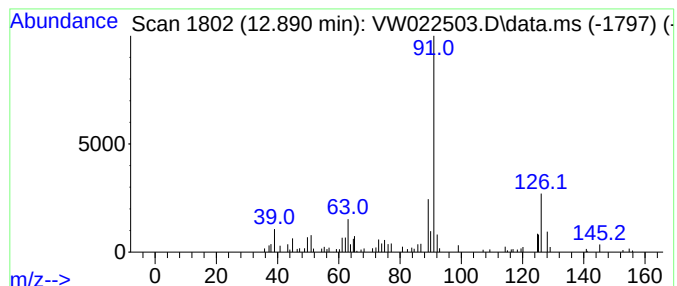
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM040722SMA.M
Quant Title : SFAM01.0

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

Peak Number 7 Benzene, 1-chloro-4-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.890	5.80 ug/L	33561	1,4-Dichlorobenzene-d4	13.554

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-chloro-4-methyl-	126	C7H7Cl	000106-43-4	92
2			Benzene, 1-chloro-2-methyl-	126	C7H7Cl	000095-49-8	81
3			Benzene, 1-chloro-2-methyl-	126	C7H7Cl	000095-49-8	76
4			Benzene, 1-chloro-3-methyl-	126	C7H7Cl	000108-41-8	70
5			Benzene, 1-chloro-2-methyl-	126	C7H7Cl	000095-49-8	68



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW041222\
Data File : VW022503.D
Acq On : 12 Apr 2022 13:33
Operator : SY/VA
Sample : N2316-12RE
Misc : 7.42g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
ETNR2RE

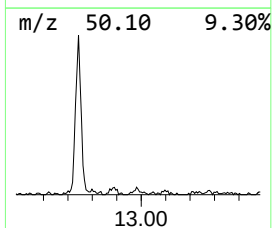
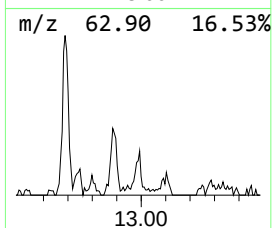
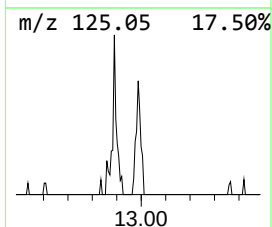
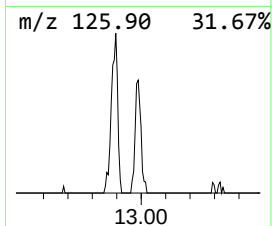
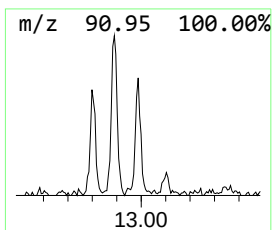
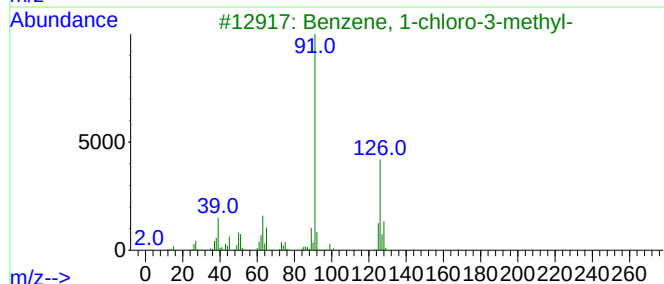
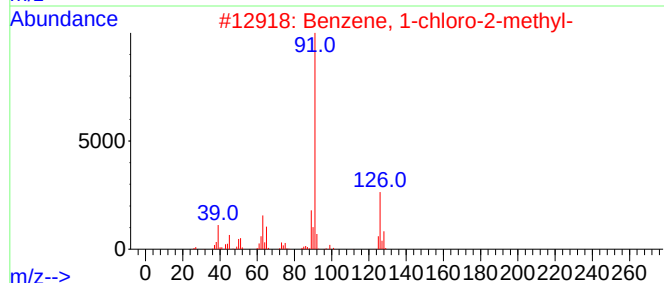
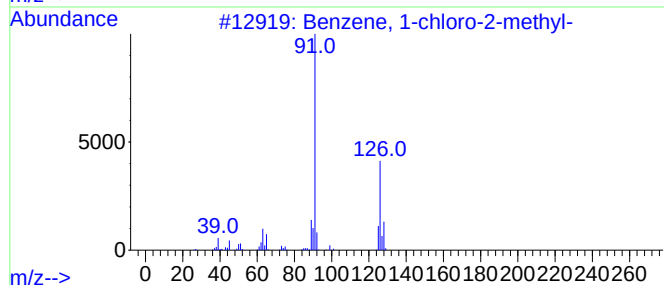
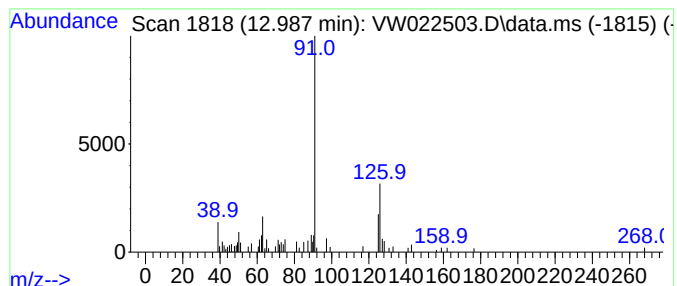
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM040722SMA.M
Quant Title : SFAM01.0

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

Peak Number 8 Benzene, 1-chloro-2-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.987	2.71 ug/L	15675	1,4-Dichlorobenzene-d4	13.554

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-chloro-2-methyl-	126	C7H7Cl	000095-49-8	58
2			Benzene, 1-chloro-2-methyl-	126	C7H7Cl	000095-49-8	53
3			Benzene, 1-chloro-3-methyl-	126	C7H7Cl	000108-41-8	52
4			Benzene, 1-chloro-4-methyl-	126	C7H7Cl	000106-43-4	49
5			Benzene, 1-chloro-4-methyl-	126	C7H7Cl	000106-43-4	49



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW041222\
Data File : VW022503.D
Acq On : 12 Apr 2022 13:33
Operator : SY/VA
Sample : N2316-12RE
Misc : 7.42g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
ETNR2RE

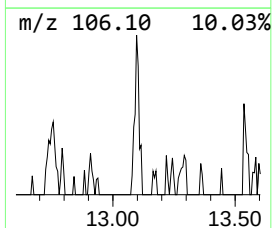
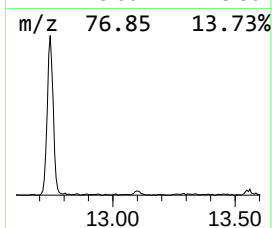
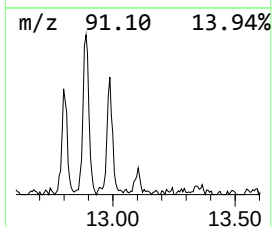
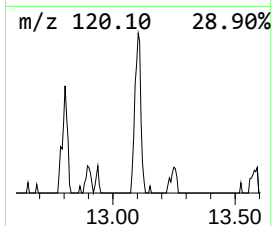
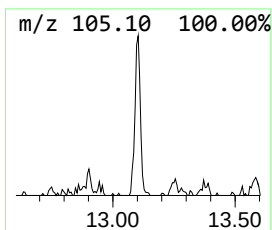
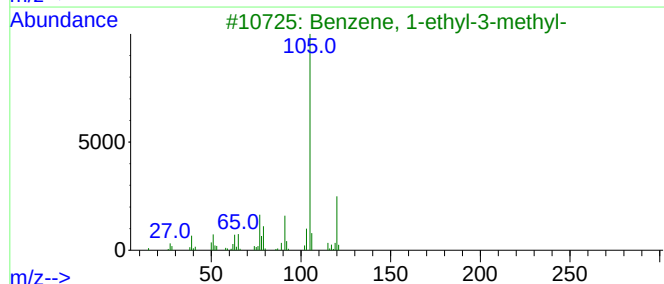
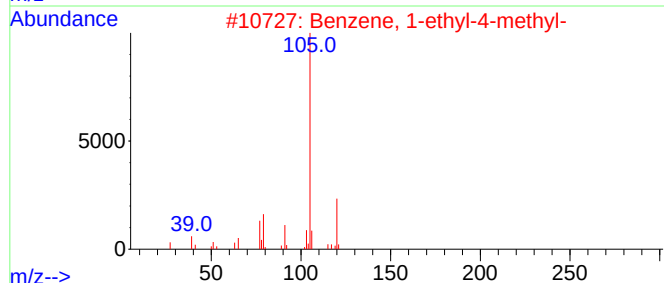
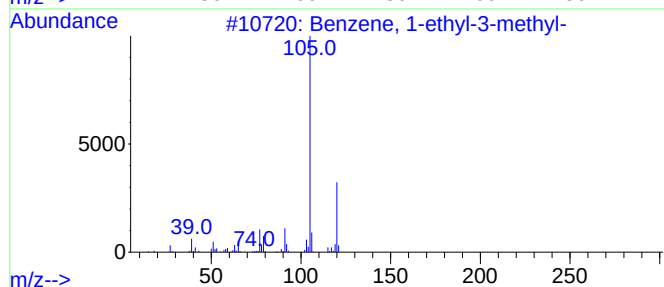
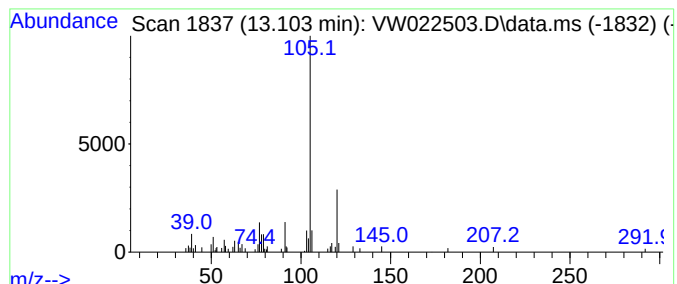
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM040722SMA.M
Quant Title : SFAM01.0

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

Peak Number 9 Benzene, 1-ethyl-3-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.103	3.87 ug/L	22424	1,4-Dichlorobenzene-d4	13.554

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
2			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	90
3			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	90
4			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	90
5			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW041222\
Data File : VW022503.D
Acq On : 12 Apr 2022 13:33
Operator : SY/VA
Sample : N2316-12RE
Misc : 7.42g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
ETNR2RE

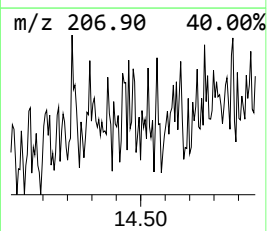
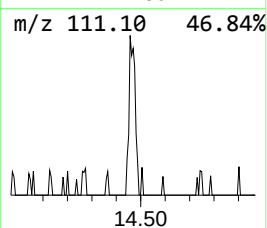
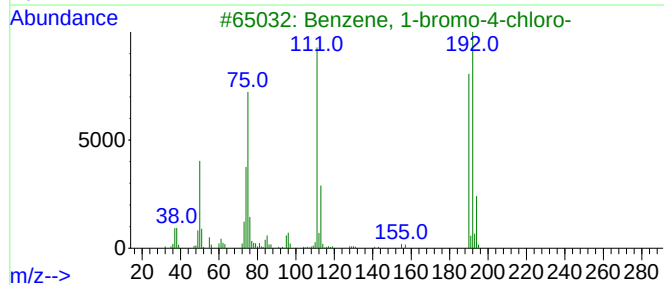
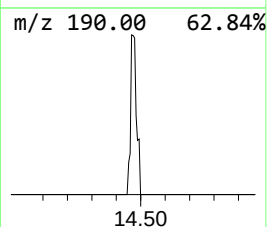
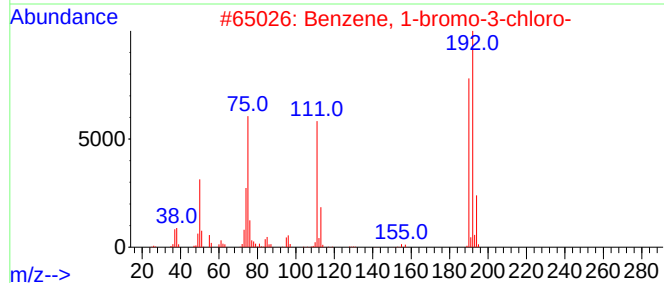
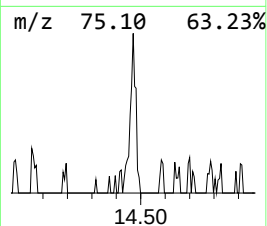
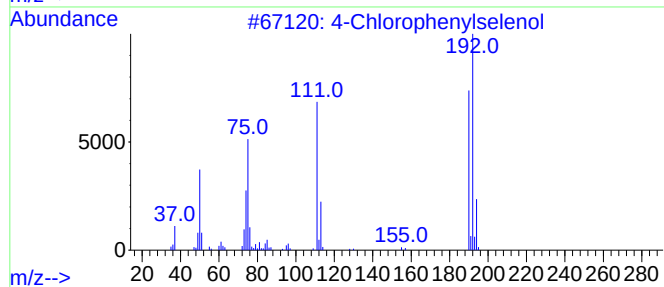
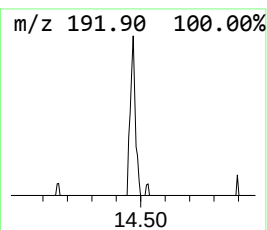
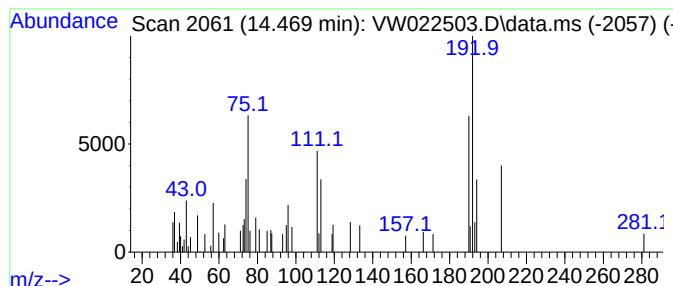
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM040722SMA.M
Quant Title : SFAM01.0

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

Peak Number 10 4-Chlorophenylselenol Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.469	1.50 ug/L	8672	1,4-Dichlorobenzene-d4	13.554

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4-Chlorophenylselenol	192	C6H5ClSe	016645-10-6	58
2			Benzene, 1-bromo-3-chloro-	190	C6H4BrCl	000108-37-2	52
3			Benzene, 1-bromo-4-chloro-	190	C6H4BrCl	000106-39-8	46
4			Benzene, 1-bromo-3-chloro-	190	C6H4BrCl	000108-37-2	46
5			Benzene, 1-bromo-2-chloro-	190	C6H4BrCl	000694-80-4	46



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW041222\
Data File : VW022503.D
Acq On : 12 Apr 2022 13:33
Operator : SY/VA
Sample : N2316-12RE
Misc : 7.42g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
ETNR2RE

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM040722SMA.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
(DEL) Alkane: S...	1.959	35.6	ug/L	176300	1	8.842	123875	25.0
unknown-01	5.574	2.3	ug/L	11265	1	8.842	123875	25.0
Thietane	9.512	3.1	ug/L	15480	1	8.842	123875	25.0
2H-Pyran, tetra...	9.957	4.0	ug/L	19583	1	8.842	123875	25.0
Benzene, bromo-	12.743	27.4	ug/L	158467	3	13.554	144746	25.0
Benzene, 1-chlo...	12.890	5.8	ug/L	33561	3	13.554	144746	25.0
Benzene, 1-chlo...	12.987	2.7	ug/L	15675	3	13.554	144746	25.0
Benzene, 1-ethy...	13.103	3.9	ug/L	22424	3	13.554	144746	25.0
4-Chlorophenyls...	14.469	1.5	ug/L	8672	3	13.554	144746	25.0