

Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW020222\
 Data File : VW022026.D
 Acq On : 02 Feb 2022 13:03
 Operator : SY/VA
 Sample : N1353-08
 Misc : 3.91g/10mL/MSVOA_W/SOIL
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 C0VF1

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\SFAMWLM020122SMA.M
 Title : SFAM01.0

Signal : TIC: VW022026.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.154	38	41	44	rBV2	660	908	0.03%	0.008%
2	2.349	67	73	81	rVB	44513	71078	2.19%	0.599%
3	2.879	155	160	173	rVB	26959	55903	1.72%	0.471%
4	3.294	227	228	230	rBV	225	134	0.00%	0.001%
5	3.367	237	240	243	rBV3	488	728	0.02%	0.006%
6	3.495	258	261	262	rBV	335	233	0.01%	0.002%
7	3.525	265	266	269	rVB	209	143	0.00%	0.001%
8	3.611	276	280	281	rBV	293	222	0.01%	0.002%
9	3.653	284	287	289	rVB2	328	315	0.01%	0.003%
10	3.678	289	291	294	rVB	220	239	0.01%	0.002%
11	3.708	294	296	297	rBV	193	200	0.01%	0.002%
12	3.849	317	319	323	rBV	194	258	0.01%	0.002%
13	4.019	336	347	358	rBV2	50098	143069	4.41%	1.206%
14	4.141	358	367	381	rVV	29647	90853	2.80%	0.766%
15	4.464	410	420	421	rBV	3168	7023	0.22%	0.059%
16	4.653	449	451	452	rBV	418	317	0.01%	0.003%
17	4.739	462	465	468	rBV	335	393	0.01%	0.003%
18	4.800	473	475	477	rBV2	198	184	0.01%	0.002%
19	4.903	486	492	493	rBV3	1106	1606	0.05%	0.014%
20	4.989	504	506	508	rVB2	181	129	0.00%	0.001%
21	5.104	521	525	526	rBV2	591	764	0.02%	0.006%
22	5.397	571	573	577	rBV	160	215	0.01%	0.002%
23	5.434	577	579	581	rBV	213	234	0.01%	0.002%
24	5.501	588	590	594	rVB	332	299	0.01%	0.003%
25	5.562	596	600	603	rBV	179	266	0.01%	0.002%
26	5.641	612	613	616	rVB2	233	156	0.00%	0.001%
27	5.775	630	635	637	rBV2	457	787	0.02%	0.007%
28	5.848	645	647	651	rVB2	283	306	0.01%	0.003%
29	5.885	651	653	654	rBV	315	221	0.01%	0.002%
30	5.946	659	663	668	rVB3	490	868	0.03%	0.007%
31	5.988	668	670	674	rBV2	161	228	0.01%	0.002%
32	6.068	679	683	686	rBV2	226	329	0.01%	0.003%
33	6.135	692	694	696	rBV	261	147	0.00%	0.001%
34	6.348	727	729	732	rBV	142	153	0.00%	0.001%
35	6.440	741	744	745	rBV	189	222	0.01%	0.002%
36	6.470	748	749	751	rBV2	269	200	0.01%	0.002%

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Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM020122SMA.M
 Title : SFAM01.0

37	6.537	758	760	763	rVB	230	153	0.00%	0.001%
38	6.568	763	765	767	rBV2	161	157	0.00%	0.001%
39	6.604	769	771	773	rVB	336	267	0.01%	0.002%
40	6.629	773	775	776	rBV	198	149	0.00%	0.001%
41	6.665	778	781	783	rBV	296	362	0.01%	0.003%
42	7.013	836	838	839	rVB2	273	189	0.01%	0.002%
43	7.086	839	850	855	rBV	8871	24404	0.75%	0.206%
44	7.391	897	900	902	rBV	374	506	0.02%	0.004%
45	7.494	915	917	919	rBV	185	142	0.00%	0.001%
46	7.531	919	923	924	rBV3	568	689	0.02%	0.006%
47	7.647	932	942	953	rBV	66967	164562	5.07%	1.387%
48	8.079	1010	1013	1016	rBV2	265	424	0.01%	0.004%
49	8.159	1025	1026	1028	rVB	241	157	0.00%	0.001%
50	8.275	1034	1045	1059	rBV2	133114	394903	12.17%	3.328%
51	8.701	1112	1115	1119	rVB3	406	586	0.02%	0.005%
52	8.842	1129	1138	1150	rVV	154228	307917	9.49%	2.595%
53	8.957	1156	1157	1160	rBV	166	148	0.00%	0.001%
54	8.988	1160	1162	1164	rVB	236	190	0.01%	0.002%
55	9.043	1166	1171	1177	rBV6	1931	4943	0.15%	0.042%
56	9.274	1202	1209	1224	rVB2	100855	215524	6.64%	1.816%
57	9.610	1260	1264	1267	rVB3	1106	1287	0.04%	0.011%
58	9.756	1283	1288	1293	rBV5	859	1754	0.05%	0.015%
59	9.817	1297	1298	1301	rBV	313	292	0.01%	0.002%
60	9.896	1305	1311	1315	rBV5	1934	3584	0.11%	0.030%
61	9.957	1316	1321	1327	rBV3	3708	8057	0.25%	0.068%
62	10.030	1328	1333	1339	rBV	43257	75594	2.33%	0.637%
63	10.207	1355	1362	1372	rBV	69443	126242	3.89%	1.064%
64	10.323	1374	1381	1386	rVV	217472	381574	11.76%	3.215%
65	10.384	1386	1391	1399	rVB	81980	149162	4.60%	1.257%
66	10.500	1403	1410	1416	rVB2	15624	28346	0.87%	0.239%
67	10.573	1416	1422	1431	rBV	31434	58789	1.81%	0.495%
68	10.664	1432	1437	1440	rBV2	9975	17381	0.54%	0.146%
69	10.762	1448	1453	1463	rVB3	11148	20258	0.62%	0.171%
70	10.847	1463	1467	1472	rVB2	3682	5631	0.17%	0.047%
71	10.920	1472	1479	1487	rBV2	42122	79904	2.46%	0.673%
72	11.109	1506	1510	1513	rBV3	9602	15260	0.47%	0.129%
73	11.225	1525	1529	1535	rVB2	16249	28128	0.87%	0.237%
74	11.335	1543	1547	1551	rBV2	8400	12972	0.40%	0.109%
75	11.512	1571	1576	1582	rVB	16408	27027	0.83%	0.228%
76	11.628	1588	1595	1600	rBV	203065	361415	11.14%	3.045%

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 Title : SFAM01.0

77	11.731	1606	1612	1621	rVB	357182	603647	18.61%	5.086%
78	11.835	1621	1629	1639	rBV	1760693	3243617	100.00%	27.331%
79	11.969	1648	1651	1654	rBV4	2243	3530	0.11%	0.030%
80	12.164	1672	1683	1689	rBV	1045657	1759788	54.25%	14.828%
81	12.231	1689	1694	1701	rVB2	88467	162074	5.00%	1.366%
82	12.359	1702	1715	1722	rBV7	27433	115193	3.55%	0.971%
83	12.457	1727	1731	1736	rBV2	35752	60673	1.87%	0.511%
84	12.603	1752	1755	1759	rVB	37079	48987	1.51%	0.413%
85	12.688	1765	1769	1775	rVB	67953	111530	3.44%	0.940%
86	12.865	1792	1798	1801	rBV	184545	306586	9.45%	2.583%
87	12.932	1805	1809	1815	rVB2	301389	575537	17.74%	4.850%
88	13.249	1856	1861	1865	rBV	316964	481708	14.85%	4.059%
89	13.554	1907	1911	1914	rBV2	225813	367264	11.32%	3.095%
90	13.865	1956	1962	1968	rVB2	377242	708632	21.85%	5.971%
91	14.719	2098	2102	2107	rVB2	108431	153900	4.74%	1.297%
92	14.786	2109	2113	2120	rBV2	82780	169481	5.23%	1.428%
93	15.359	2202	2207	2213	rVB	43622	74705	2.30%	0.629%
94	15.432	2215	2219	2223	rVB2	7067	10850	0.33%	0.091%
95	15.651	2251	2255	2263	rVB6	1539	3093	0.10%	0.026%
96	15.944	2296	2303	2309	rBV6	1032	2431	0.07%	0.020%
97	16.218	2346	2348	2350	rBV	424	274	0.01%	0.002%
98	16.322	2363	2365	2366	rBV	480	272	0.01%	0.002%
99	16.432	2378	2383	2389	rBV4	1960	3903	0.12%	0.033%
100	16.633	2412	2416	2423	rVB4	936	1670	0.05%	0.014%

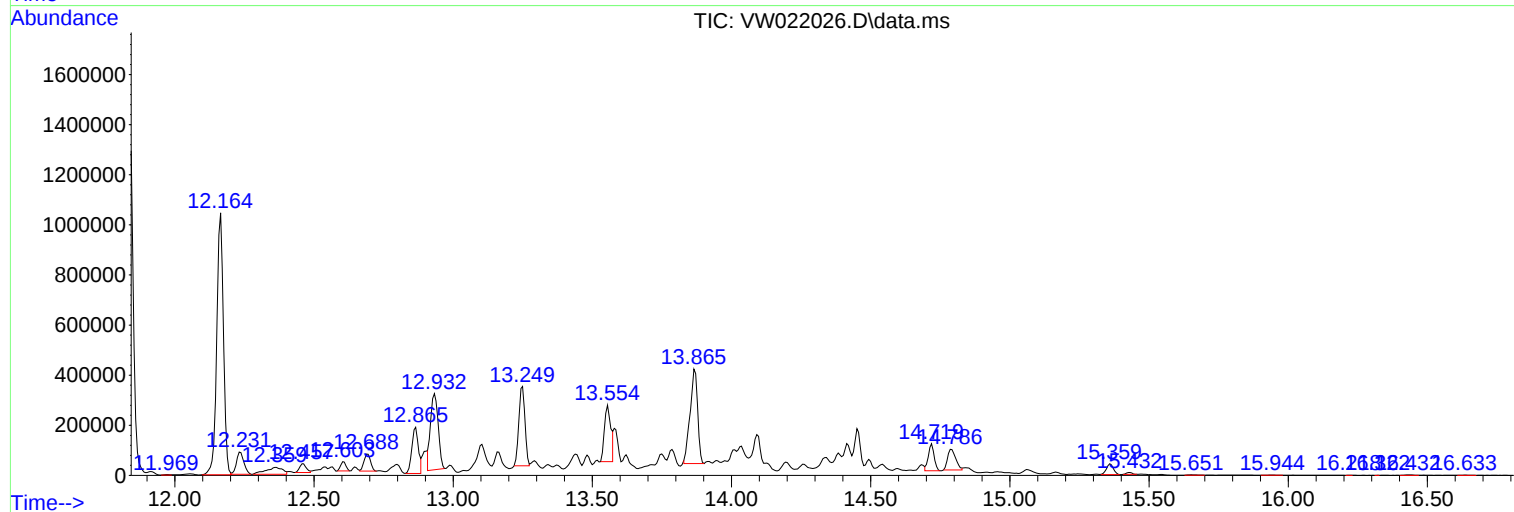
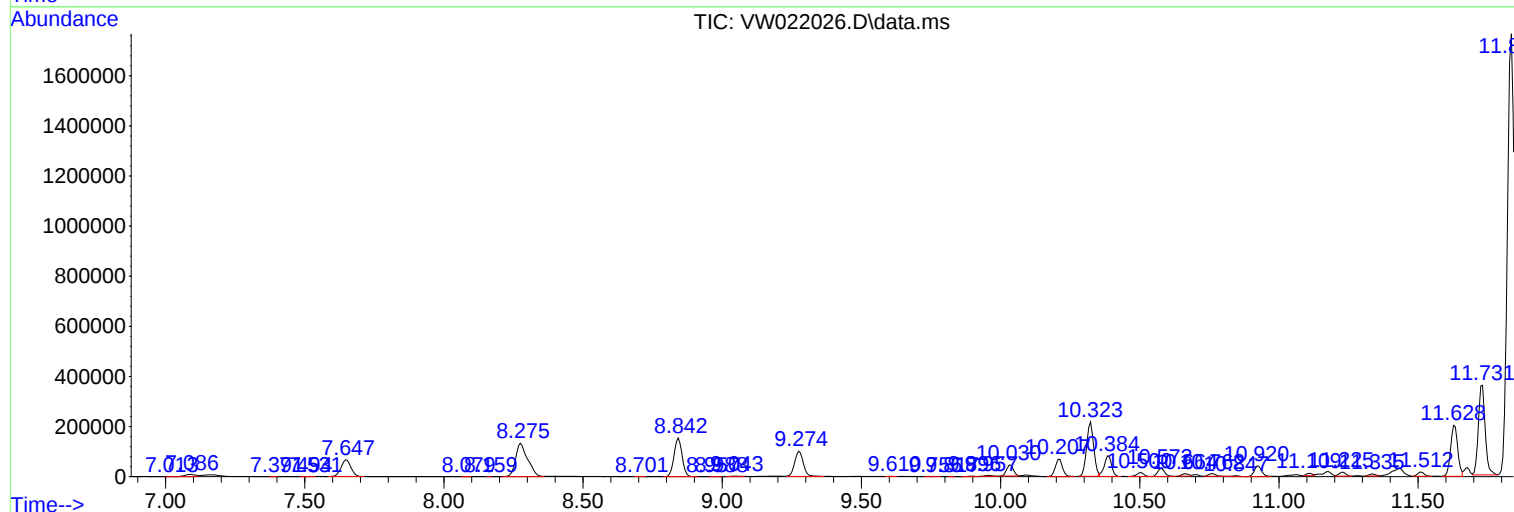
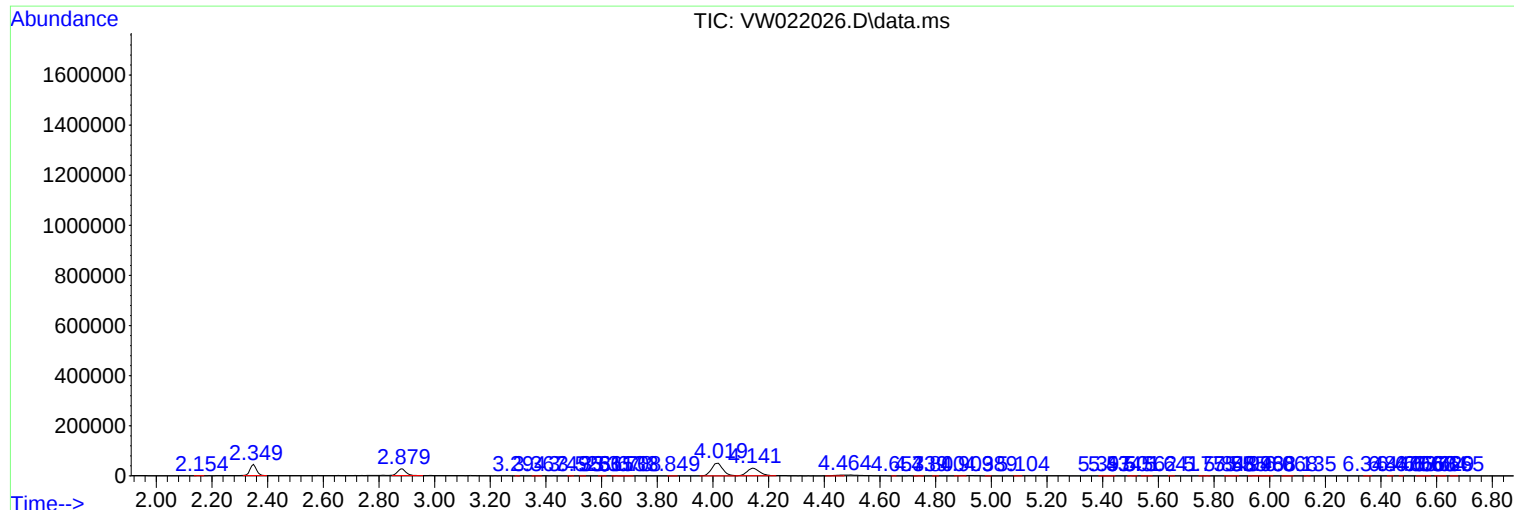
Sum of corrected areas: 11867704

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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM020122SMA.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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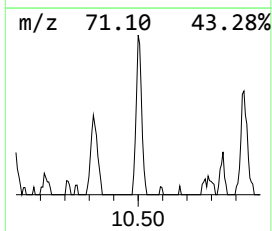
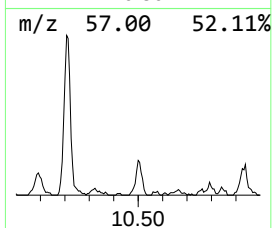
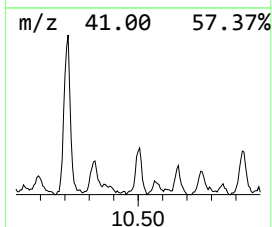
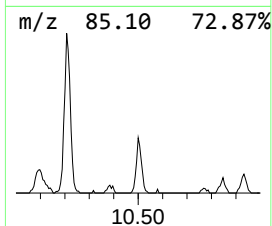
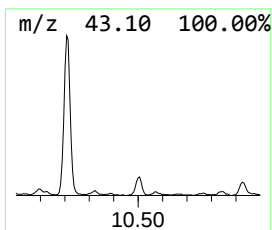
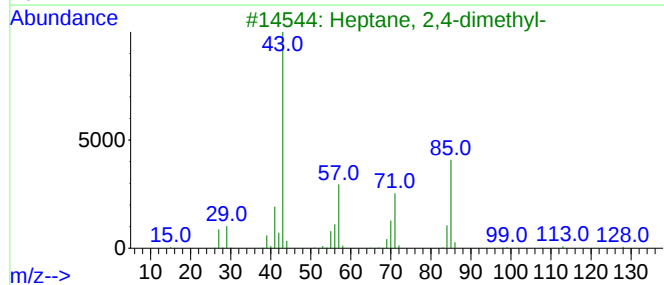
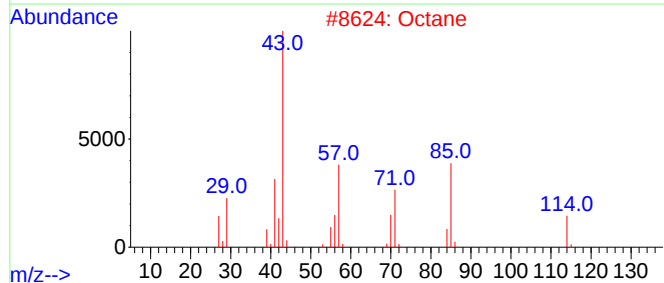
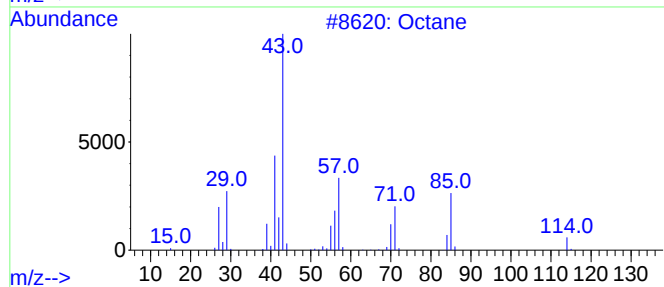
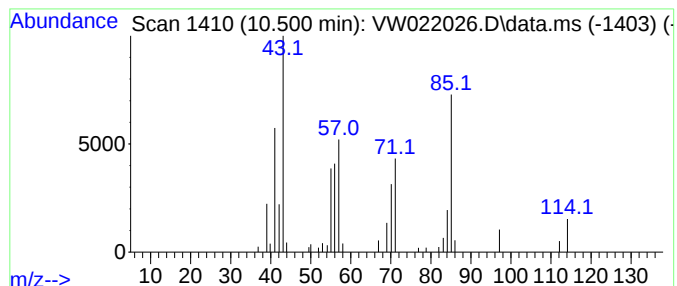
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM020122SMA.M
Quant Title : SFAM01.0

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

Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.500	1.96 ug/L	28346	Chlorobenzene-d5	11.634

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octane	114	C8H18	000111-65-9	87
2			Octane	114	C8H18	000111-65-9	64
3			Heptane, 2,4-dimethyl-	128	C9H20	002213-23-2	50
4			Hexane, 3,3-dimethyl-	114	C8H18	000563-16-6	47
5			1-Octene, 4-methyl-	126	C9H18	013151-12-7	43



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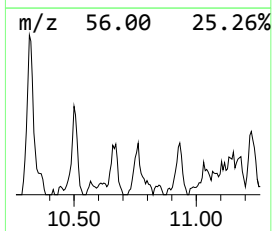
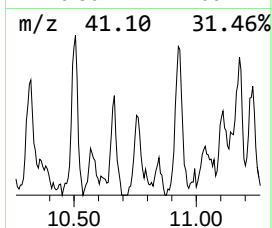
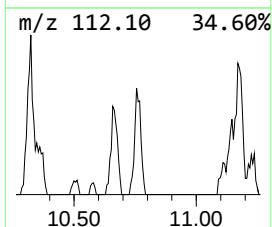
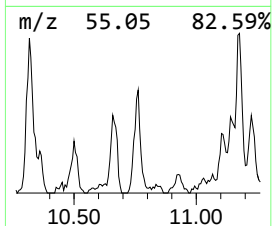
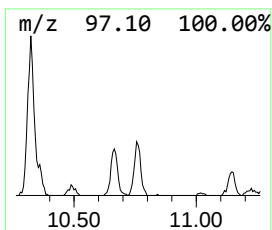
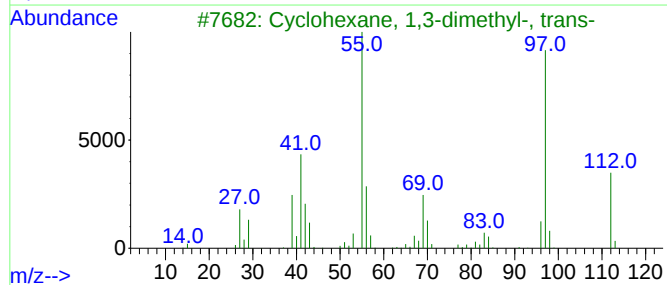
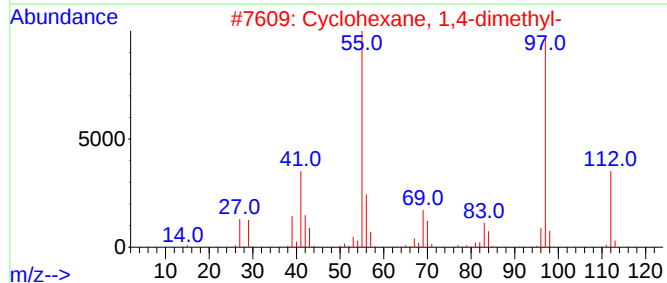
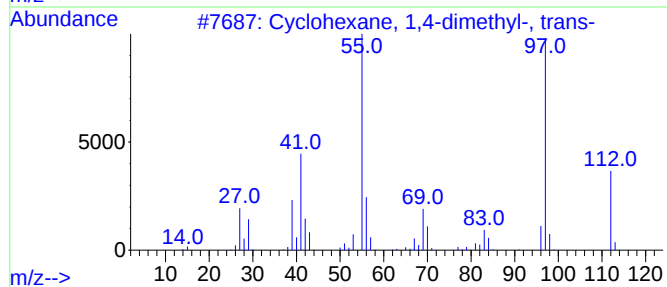
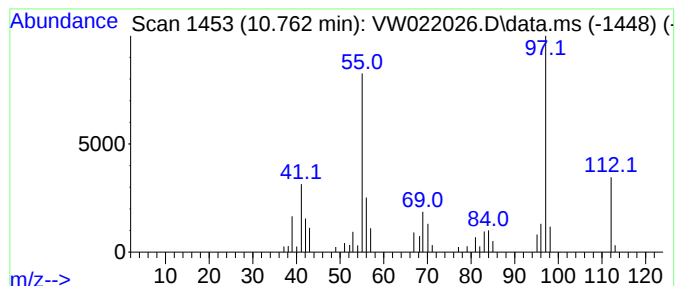
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM020122SMA.M
Quant Title : SFAM01.0

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

Peak Number 3 (DEL) Alkane: Cyclic10.762 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.762	1.40 ug/L	20258	Chlorobenzene-d5	11.634

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, 1,4-dimethyl-, trans-	112	C8H16	002207-04-7	97
2			Cyclohexane, 1,4-dimethyl-	112	C8H16	000589-90-2	95
3			Cyclohexane, 1,3-dimethyl-, trans-	112	C8H16	002207-03-6	94
4			Cyclohexane, 1,4-dimethyl-, cis-	112	C8H16	000624-29-3	93
5			Cyclohexane, 1,4-dimethyl-, cis-	112	C8H16	000624-29-3	93



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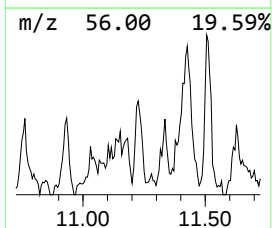
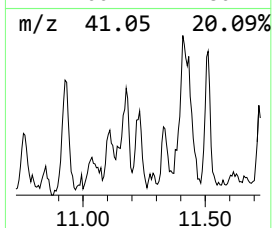
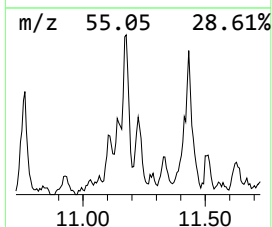
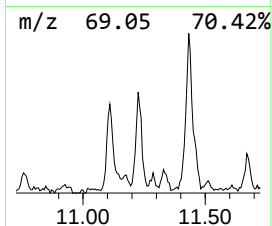
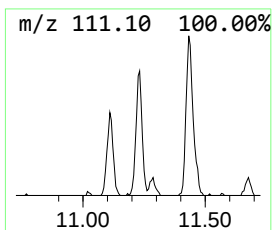
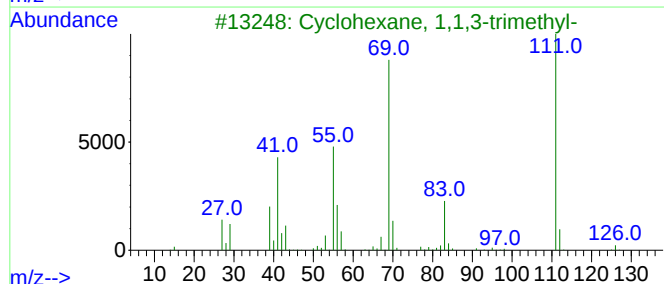
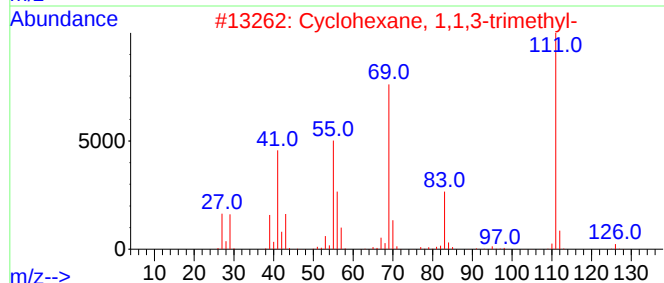
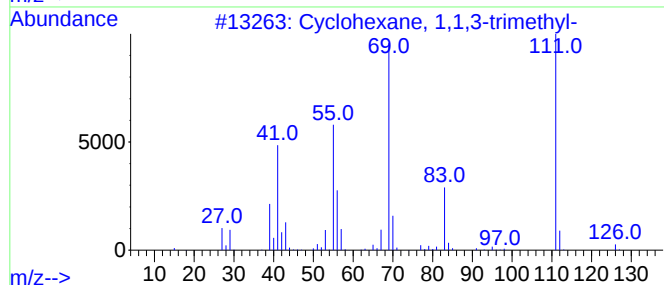
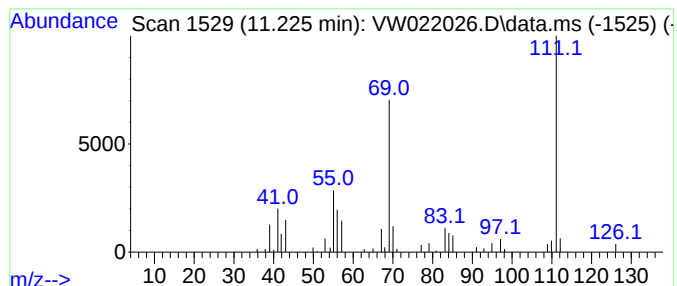
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM020122SMA.M
Quant Title : SFAM01.0

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

Peak Number 4 (DEL) Alkane: Cyclic11.225 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.225	1.95 ug/L	28128	Chlorobenzene-d5	11.634

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, 1,1,3-trimethyl-	126	C9H18	003073-66-3	74
2			Cyclohexane, 1,1,3-trimethyl-	126	C9H18	003073-66-3	72
3			Cyclohexane, 1,1,3-trimethyl-	126	C9H18	003073-66-3	64
4			1,1,4-Trimethylcyclohexane	126	C9H18	007094-27-1	59
5			Pyrazol-4-amine, 1,5-dimethyl-	111	C5H9N3	121983-36-6	59



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW020222\
Data File : VW022026.D
Acq On : 02 Feb 2022 13:03
Operator : SY/VA
Sample : N1353-08
Misc : 3.91g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

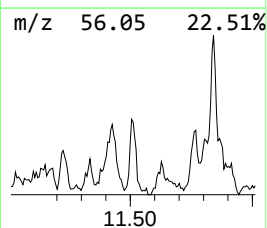
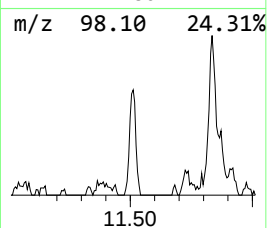
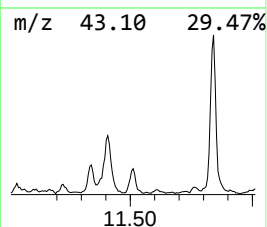
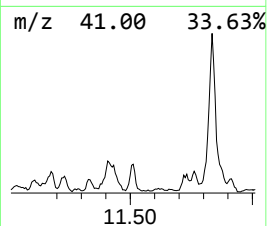
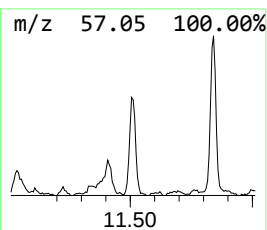
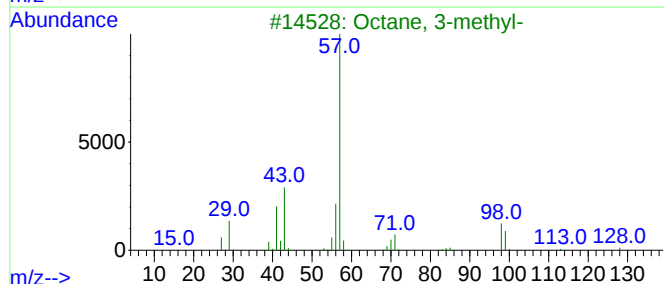
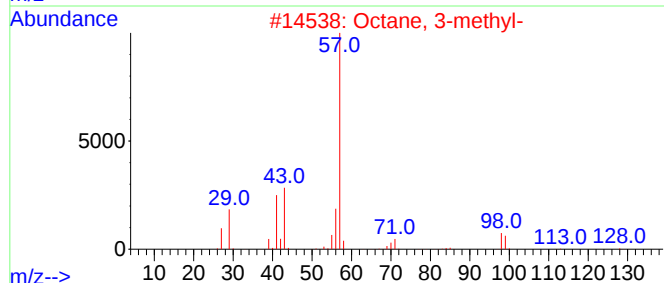
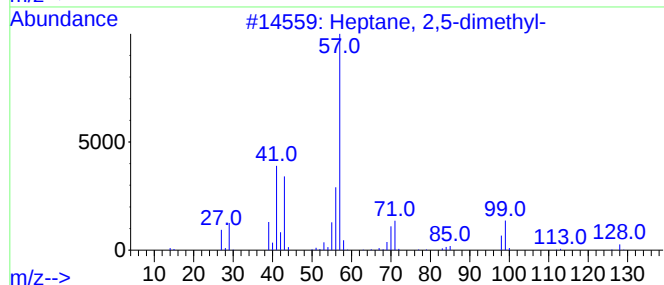
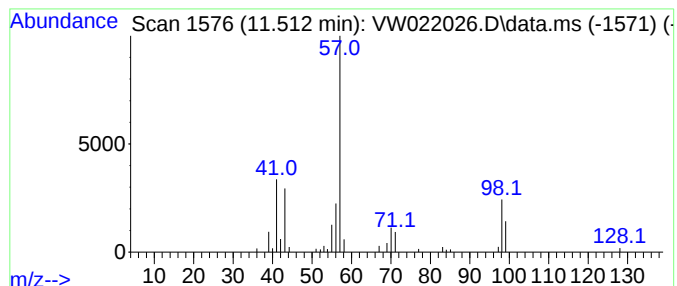
Instrument :
MSVOA_W
ClientSampleId :
COVF1

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM020122SMA.M
Quant Title : SFAM01.0

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

Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD		R.T.	
11.512	1.87 ug/L	27027	Chlorobenzene-d5		11.634	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Heptane, 2,5-dimethyl-		128	C9H20	002216-30-0	81
2	Octane, 3-methyl-		128	C9H20	002216-33-3	72
3	Octane, 3-methyl-		128	C9H20	002216-33-3	72
4	Decane, 2,5-dimethyl-		170	C12H26	017312-50-4	64
5	Undecane, 6-methyl-		170	C12H26	017302-33-9	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW020222\
Data File : VW022026.D
Acq On : 02 Feb 2022 13:03
Operator : SY/VA
Sample : N1353-08
Misc : 3.91g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
COVF1

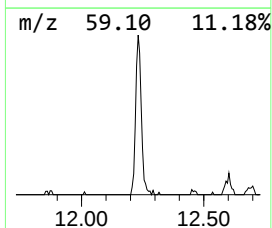
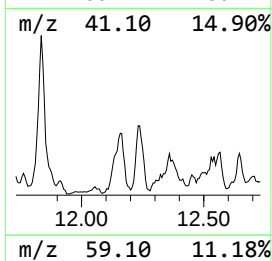
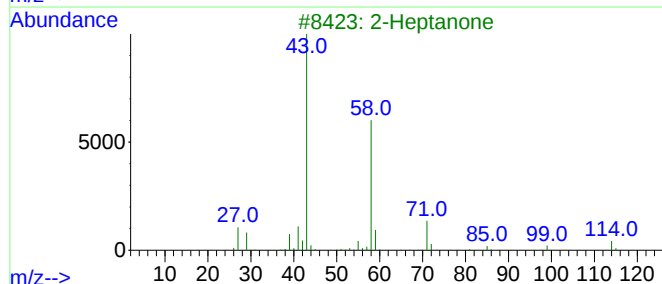
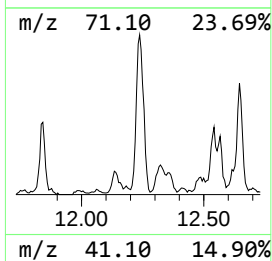
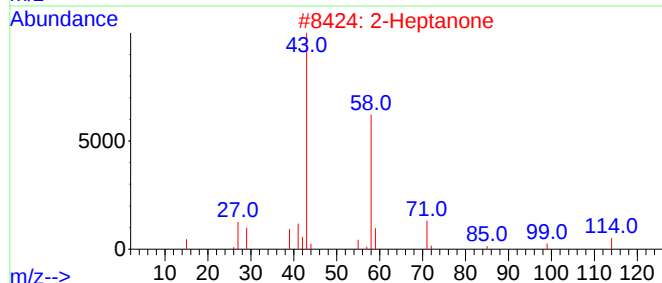
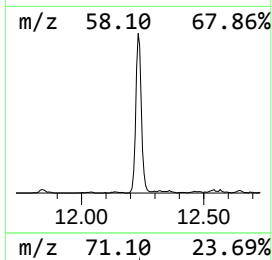
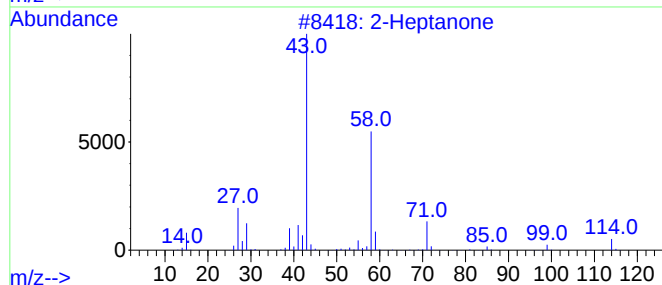
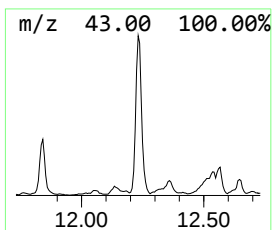
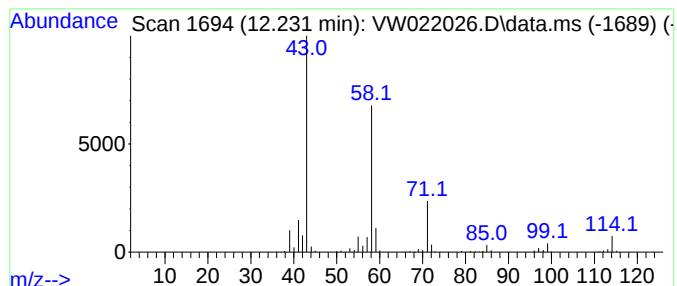
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Quant Title : SFAM01.0

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

Peak Number 6 2-Heptanone Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.231	11.21 ug/L	162074	Chlorobenzene-d5	11.634

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Heptanone			114	C7H14O	000110-43-0	91
2	2-Heptanone			114	C7H14O	000110-43-0	91
3	2-Heptanone			114	C7H14O	000110-43-0	91
4	2-Heptanone			114	C7H14O	000110-43-0	80
5	2-Heptanone			114	C7H14O	000110-43-0	80



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW020222\
Data File : VW022026.D
Acq On : 02 Feb 2022 13:03
Operator : SY/VA
Sample : N1353-08
Misc : 3.91g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

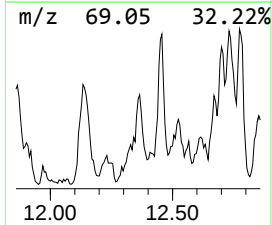
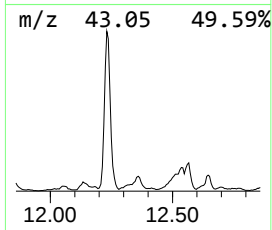
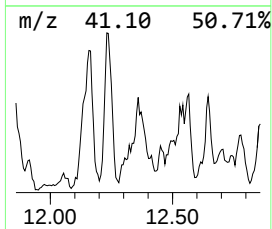
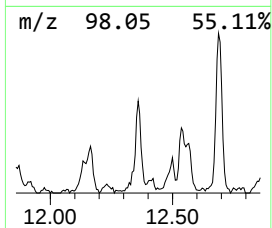
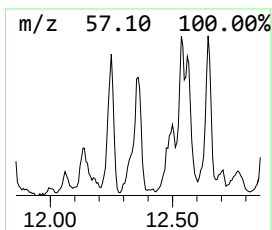
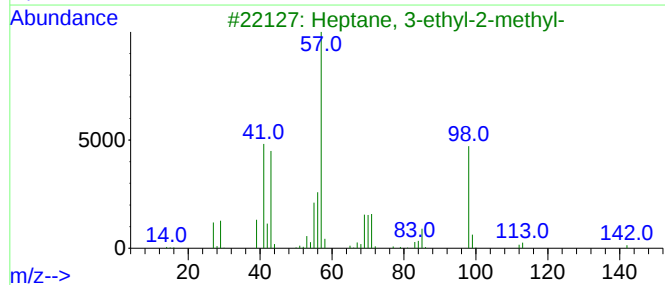
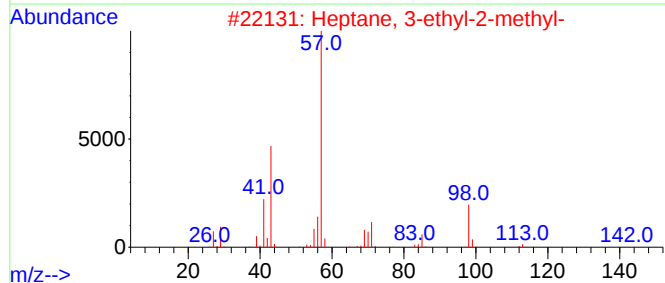
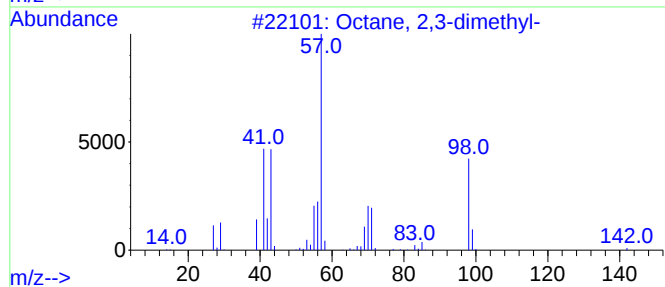
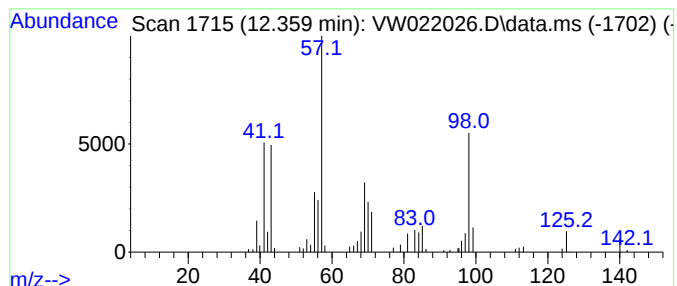
Instrument :
MSVOA_W
ClientSampleId :
COVF1

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM020122SMA.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 5

R.T.	EstConc	Area	Relative to ISTD			R.T.
12.359	7.97 ug/L	115193	Chlorobenzene-d5			11.634
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Octane, 2,3-dimethyl-		142	C10H22	007146-60-3	64
2	Heptane, 3-ethyl-2-methyl-		142	C10H22	014676-29-0	50
3	Heptane, 3-ethyl-2-methyl-		142	C10H22	014676-29-0	46
4	2,3-2H-4-Methyl-imidazole-2-one		98	C4H6N2O	039799-77-4	43
5	Heptane, 3-ethyl-2-methyl-		142	C10H22	014676-29-0	43



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW020222\
Data File : VW022026.D
Acq On : 02 Feb 2022 13:03
Operator : SY/VA
Sample : N1353-08
Misc : 3.91g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
COVF1

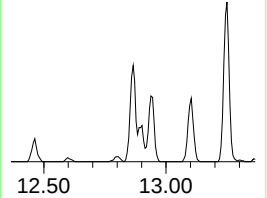
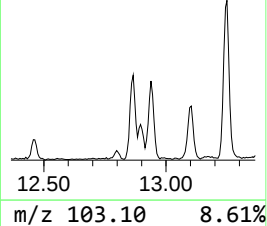
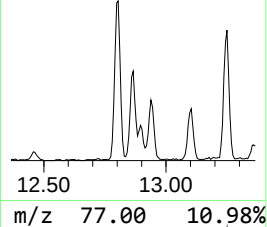
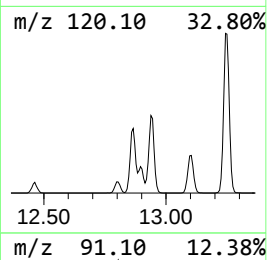
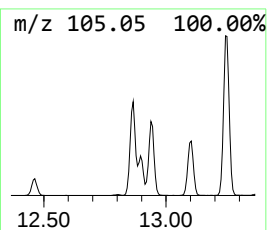
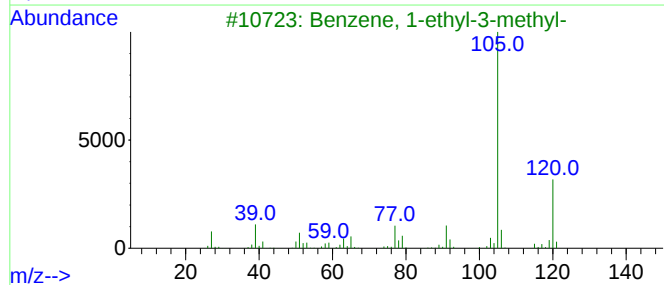
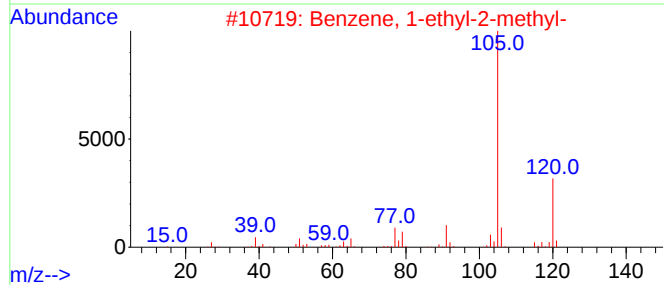
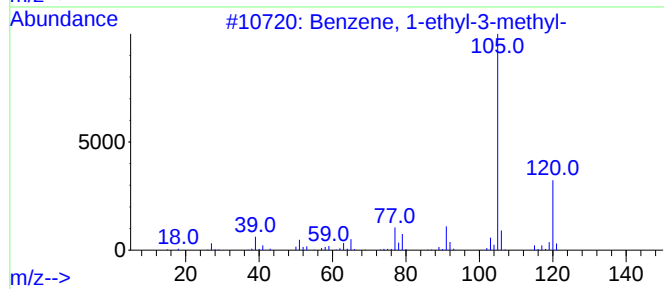
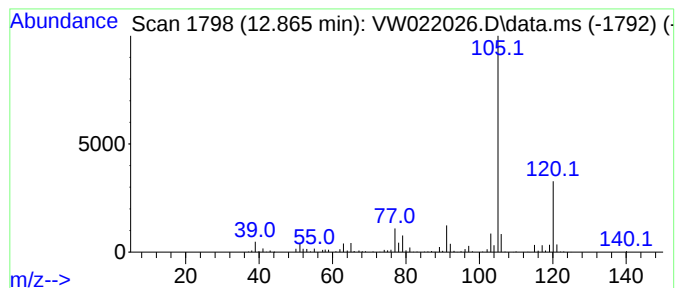
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Quant Title : SFAM01.0

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.865	20.87 ug/L	306586	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	95
2			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
3			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	94
4			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94
5			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW020222\
Data File : VW022026.D
Acq On : 02 Feb 2022 13:03
Operator : SY/VA
Sample : N1353-08
Misc : 3.91g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
COVF1

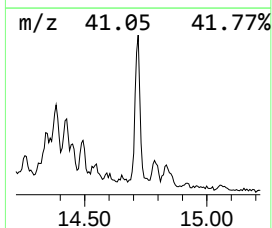
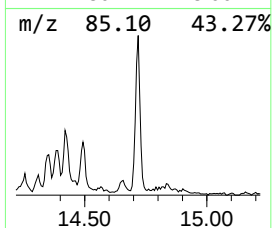
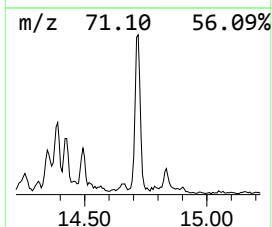
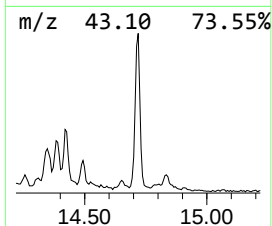
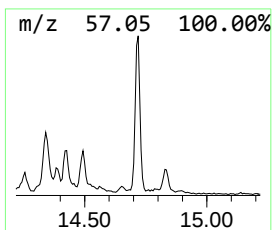
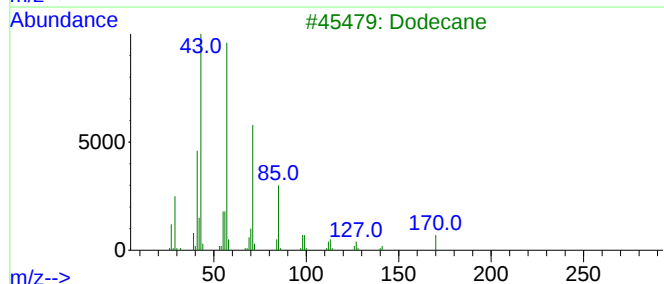
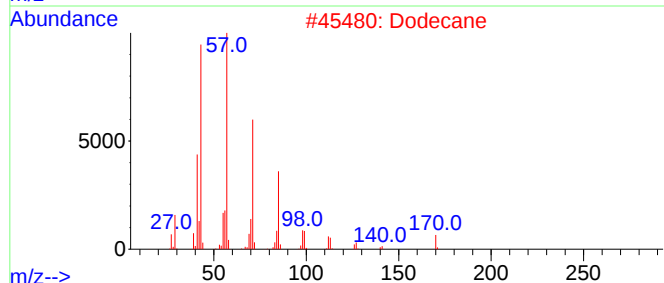
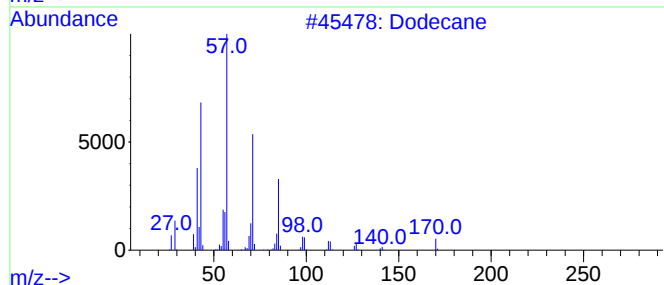
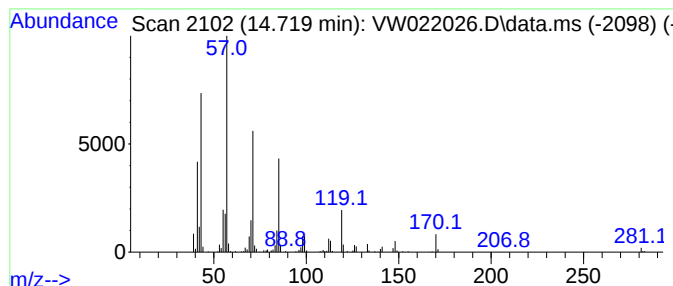
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Quant Title : SFAM01.0

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

Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.719	10.48 ug/L	153900	1,4-Dichlorobenzene-d4	13.554

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dodecane	170	C12H26	000112-40-3	96
2			Dodecane	170	C12H26	000112-40-3	94
3			Dodecane	170	C12H26	000112-40-3	89
4			Dodecane	170	C12H26	000112-40-3	76
5			Tridecane	184	C13H28	000629-50-5	62



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW020222\
Data File : VW022026.D
Acq On : 02 Feb 2022 13:03
Operator : SY/VA
Sample : N1353-08
Misc : 3.91g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
COVF1

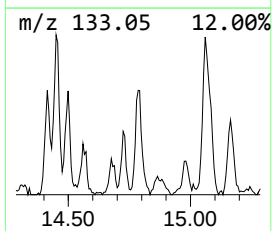
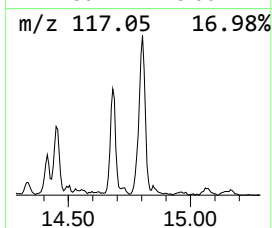
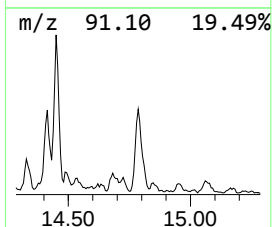
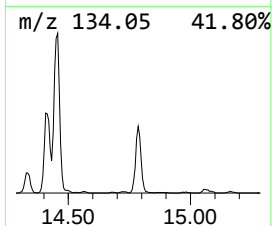
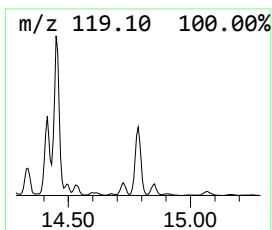
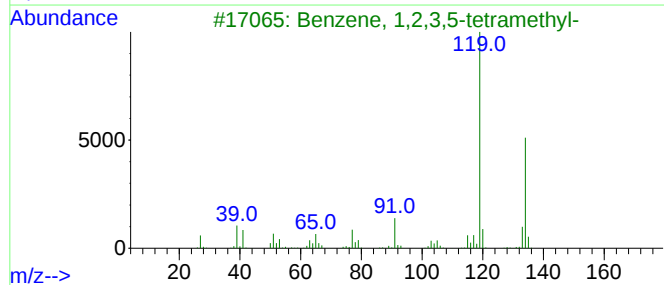
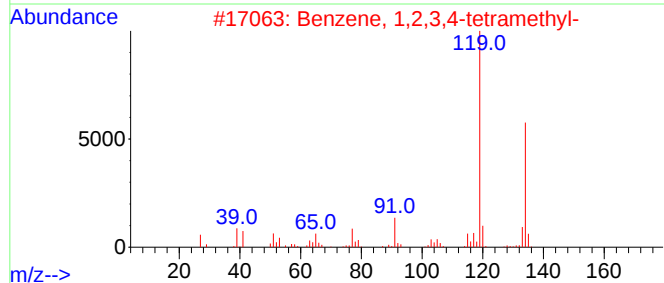
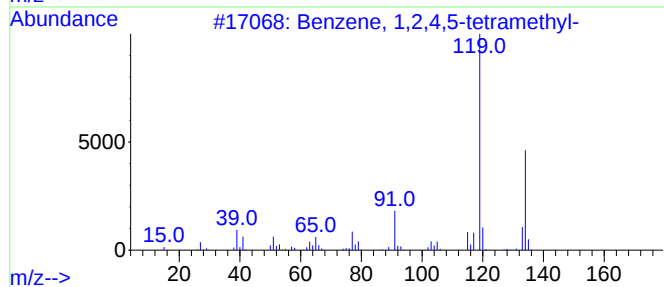
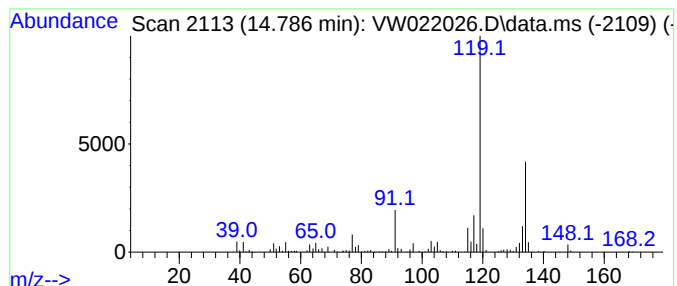
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Quant Title : SFAM01.0

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

Peak Number 11 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.786	11.54 ug/L	169481	1,4-Dichlorobenzene-d4	13.554

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	95
2			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	90
3			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	90
4			1,3-Cyclopentadiene, 1,2,3,4-tet...	134	C10H14	076089-59-3	90
5			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW020222\
Data File : VW022026.D
Acq On : 02 Feb 2022 13:03
Operator : SY/VA
Sample : N1353-08
Misc : 3.91g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
COVF1

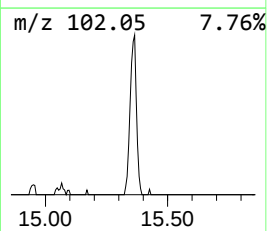
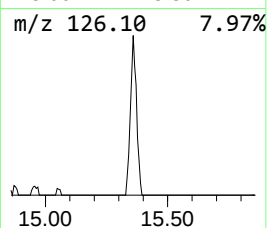
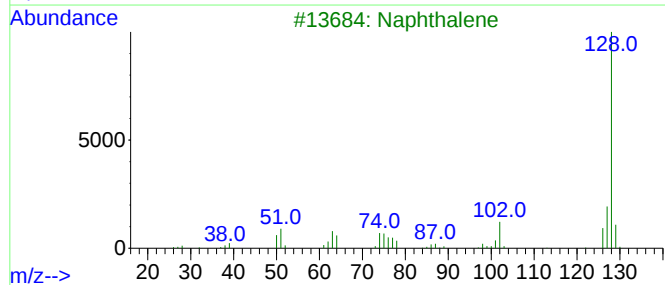
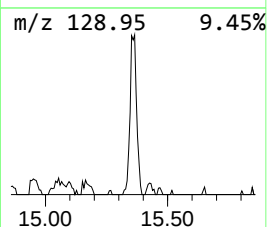
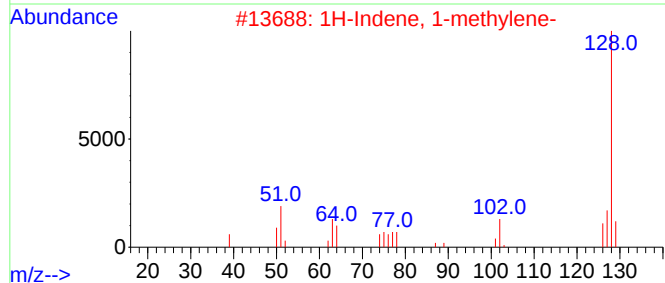
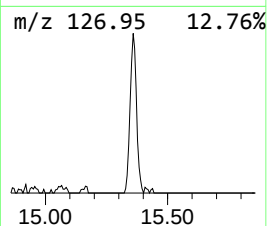
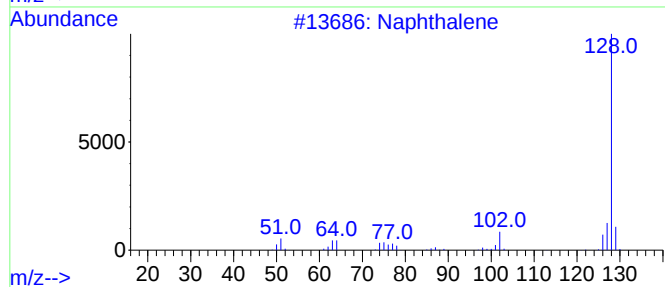
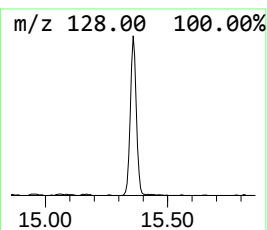
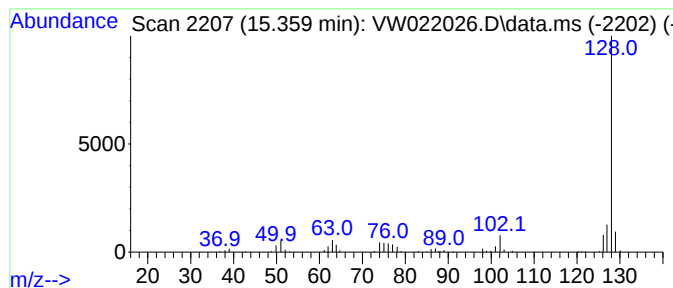
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Quant Title : SFAM01.0

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

Peak Number 12 Azulene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.359	5.09 ug/L	74705	1,4-Dichlorobenzene-d4	13.554

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene	128	C10H8	000091-20-3	94
2			1H-Indene, 1-methylene-	128	C10H8	002471-84-3	91
3			Naphthalene	128	C10H8	000091-20-3	91
4			Azulene	128	C10H8	000275-51-4	90
5			Naphthalene	128	C10H8	000091-20-3	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW020222\
Data File : VW022026.D
Acq On : 02 Feb 2022 13:03
Operator : SY/VA
Sample : N1353-08
Misc : 3.91g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
C0VF1

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM020122SMA.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	28346	2	11.634	361415	25.0
(DEL) Alkane: C...	10.762	1.4	ug/L	20258	2	11.634	361415	25.0
(DEL) Alkane: C...	11.225	2.0	ug/L	28128	2	11.634	361415	25.0
(DEL) Alkane: S...	11.512	1.9	ug/L	27027	2	11.634	361415	25.0
2-Heptanone	12.231	11.2	ug/L	162074	2	11.634	361415	25.0
(DEL) Alkane: S...	12.359	8.0	ug/L	115193	2	11.634	361415	25.0
Benzene, 1-ethy...	12.865	20.9	ug/L	306586	3	13.554	367264	25.0
(DEL) Alkane: S...	14.719	10.5	ug/L	153900	3	13.554	367264	25.0
Benzene, 1,2,4,...	14.786	11.5	ug/L	169481	3	13.554	367264	25.0
Azulene	15.359	5.1	ug/L	74705	3	13.554	367264	25.0