

Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW081221\
 Data File : VW019719.D
 Acq On : 12 Aug 2021 11:39
 Operator : SY/VA
 Sample : M3365-03
 Misc : 5.92g/10mL/MSVOA_W/SOIL
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 DBK83

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\SFAMWLM081121SMA.M
 Title : SFAM01.0

Signal : TIC: VW019719.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.361	70	75	86	rVB	71805	143990	6.08%	0.985%
2	2.891	156	162	179	rVB	73312	186019	7.85%	1.273%
3	3.507	261	263	266	rBV2	235	265	0.01%	0.002%
4	3.580	273	275	277	rVB3	345	246	0.01%	0.002%
5	3.599	277	278	281	rBV	311	242	0.01%	0.002%
6	3.666	287	289	290	rBV	292	189	0.01%	0.001%
7	3.745	299	302	305	rBV2	351	450	0.02%	0.003%
8	4.019	335	347	363	rBV	107436	327422	13.82%	2.240%
9	4.348	399	401	402	rBV	326	254	0.01%	0.002%
10	4.604	441	443	447	rVB	209	206	0.01%	0.001%
11	4.653	448	451	452	rBV2	222	218	0.01%	0.001%
12	4.672	452	454	457	rVV3	265	200	0.01%	0.001%
13	4.793	472	474	477	rVB	341	244	0.01%	0.002%
14	4.915	483	494	505	rBV	16705	60017	2.53%	0.411%
15	5.421	575	577	583	rVB2	379	638	0.03%	0.004%
16	5.580	600	603	606	rVB	164	195	0.01%	0.001%
17	5.617	606	609	611	rBV2	152	196	0.01%	0.001%
18	5.799	636	639	641	rBV2	241	190	0.01%	0.001%
19	5.946	652	663	672	rBV4	5740	17811	0.75%	0.122%
20	6.062	680	682	685	rVB	292	242	0.01%	0.002%
21	6.854	809	812	815	rBV2	227	217	0.01%	0.001%
22	6.964	827	830	832	rBV2	164	208	0.01%	0.001%
23	7.013	835	838	840	rBV2	289	328	0.01%	0.002%
24	7.086	840	850	867	rBV	30933	90112	3.80%	0.616%
25	7.513	917	920	921	rBV2	291	219	0.01%	0.001%
26	7.647	931	942	956	rBV	179394	444503	18.77%	3.041%
27	7.769	960	962	966	rVB	205	222	0.01%	0.002%
28	7.872	976	979	980	rBV3	225	222	0.01%	0.002%
29	7.945	986	991	997	rBV3	1534	3098	0.13%	0.021%
30	8.067	1008	1011	1018	rVV4	457	801	0.03%	0.005%
31	8.275	1033	1045	1062	rBV2	350131	1038710	43.85%	7.106%
32	8.585	1094	1096	1099	rBV2	227	248	0.01%	0.002%
33	8.628	1099	1103	1105	rBV2	254	299	0.01%	0.002%
34	8.842	1130	1138	1152	rBV	364548	720334	30.41%	4.928%
35	9.006	1163	1165	1166	rBV	258	236	0.01%	0.002%
36	9.049	1170	1172	1173	rVV	256	239	0.01%	0.002%

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Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM081121SMA.M
 Title : SFAM01.0

37	9.085	1173	1178	1183	rVV4	1821	3447	0.15%	0.024%
38	9.274	1198	1209	1219	rBV	243292	494262	20.87%	3.381%
39	9.512	1243	1248	1252	rVB3	228	407	0.02%	0.003%
40	9.658	1266	1272	1277	rVV3	1673	3479	0.15%	0.024%
41	10.036	1324	1334	1343	rBV	124565	228064	9.63%	1.560%
42	10.122	1344	1348	1355	rVB3	1329	2381	0.10%	0.016%
43	10.207	1357	1362	1368	rBV4	546	1164	0.05%	0.008%
44	10.323	1373	1381	1387	rBV	484549	841393	35.52%	5.756%
45	10.390	1387	1392	1399	rVB	90648	159387	6.73%	1.090%
46	10.488	1406	1408	1410	rBV	222	213	0.01%	0.001%
47	10.579	1416	1423	1434	rVB	84012	144602	6.11%	0.989%
48	10.707	1437	1444	1450	rVB2	29373	53888	2.28%	0.369%
49	10.756	1450	1452	1453	rBV2	283	223	0.01%	0.002%
50	10.859	1465	1469	1473	rBV3	1091	1384	0.06%	0.009%
51	10.927	1473	1480	1493	rVV	120556	206766	8.73%	1.414%
52	11.061	1500	1502	1508	rVB3	498	759	0.03%	0.005%
53	11.183	1517	1522	1526	rBV5	816	1702	0.07%	0.012%
54	11.274	1534	1537	1539	rVB2	236	274	0.01%	0.002%
55	11.341	1546	1548	1551	rBV3	197	226	0.01%	0.002%
56	11.408	1555	1559	1562	rVB2	228	350	0.01%	0.002%
57	11.634	1587	1596	1606	rBV	527935	877234	37.04%	6.001%
58	11.731	1608	1612	1618	rVV2	2304	3876	0.16%	0.027%
59	11.841	1626	1630	1638	rVB4	5314	9705	0.41%	0.066%
60	12.054	1664	1665	1667	rBV	287	181	0.01%	0.001%
61	12.170	1679	1684	1691	rVB3	3658	7101	0.30%	0.049%
62	12.225	1691	1693	1695	rBV2	314	309	0.01%	0.002%
63	12.463	1727	1732	1738	rBV	6218	9643	0.41%	0.066%
64	12.609	1746	1756	1763	rBV	23720	41768	1.76%	0.286%
65	12.695	1763	1770	1777	rBV	196920	328401	13.86%	2.247%
66	12.798	1784	1787	1794	rVB	2909	4729	0.20%	0.032%
67	12.871	1794	1799	1800	rBV2	855	1177	0.05%	0.008%
68	12.896	1800	1803	1807	rVV2	3375	5510	0.23%	0.038%
69	12.944	1807	1811	1817	rVB2	5629	9268	0.39%	0.063%
70	13.005	1819	1821	1824	rVB3	415	372	0.02%	0.003%
71	13.103	1829	1837	1843	rVB	7394	11289	0.48%	0.077%
72	13.188	1846	1851	1856	rBV4	1310	2494	0.11%	0.017%
73	13.249	1856	1861	1870	rVB2	6872	11823	0.50%	0.081%
74	13.377	1878	1882	1884	rBV3	929	1369	0.06%	0.009%
75	13.438	1891	1892	1895	rBV3	604	495	0.02%	0.003%
76	13.499	1898	1902	1905	rBV5	3050	4647	0.20%	0.032%

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77	13.560	1905	1912	1924	rVB	524483	868267	36.66%	5.940%
78	13.761	1933	1945	1953	rBV	231560	406477	17.16%	2.781%
79	13.853	1953	1960	1971	rVB	434542	726175	30.66%	4.968%
80	14.005	1982	1985	1990	rBV3	2904	4491	0.19%	0.031%
81	14.066	1990	1995	1996	rBV	12440	15564	0.66%	0.106%
82	14.158	2005	2010	2014	rBV	31028	50628	2.14%	0.346%
83	14.206	2014	2018	2030	rVB	73992	115219	4.86%	0.788%
84	14.395	2043	2049	2050	rBV2	14309	22572	0.95%	0.154%
85	14.487	2056	2064	2073	rBV3	20927	52682	2.22%	0.360%
86	14.627	2083	2087	2091	rBV6	3218	5874	0.25%	0.040%
87	14.688	2091	2097	2108	rVB	201857	323650	13.66%	2.214%
88	14.804	2108	2116	2122	rBV	308269	522917	22.08%	3.577%
89	14.871	2122	2127	2133	rVB2	21471	38240	1.61%	0.262%
90	14.956	2135	2141	2147	rVB3	20763	37384	1.58%	0.256%
91	15.060	2149	2158	2162	rBV3	14656	34902	1.47%	0.239%
92	15.279	2190	2194	2197	rBV4	4963	8629	0.36%	0.059%
93	15.365	2203	2208	2214	rBV	9614	17534	0.74%	0.120%
94	15.651	2249	2255	2261	rBV2	14002	26828	1.13%	0.184%
95	15.950	2299	2304	2305	rBV4	5421	8888	0.38%	0.061%
96	16.169	2335	2340	2350	rBV5	5258	12647	0.53%	0.087%
97	16.371	2362	2373	2377	rBV2	37600	90345	3.81%	0.618%
98	16.438	2377	2384	2397	rVV	1086151	2276630	96.12%	15.574%
99	16.560	2397	2404	2409	rVV	28393	67708	2.86%	0.463%
100	16.639	2409	2417	2431	rVB	1090612	2368576	100.00%	16.203%

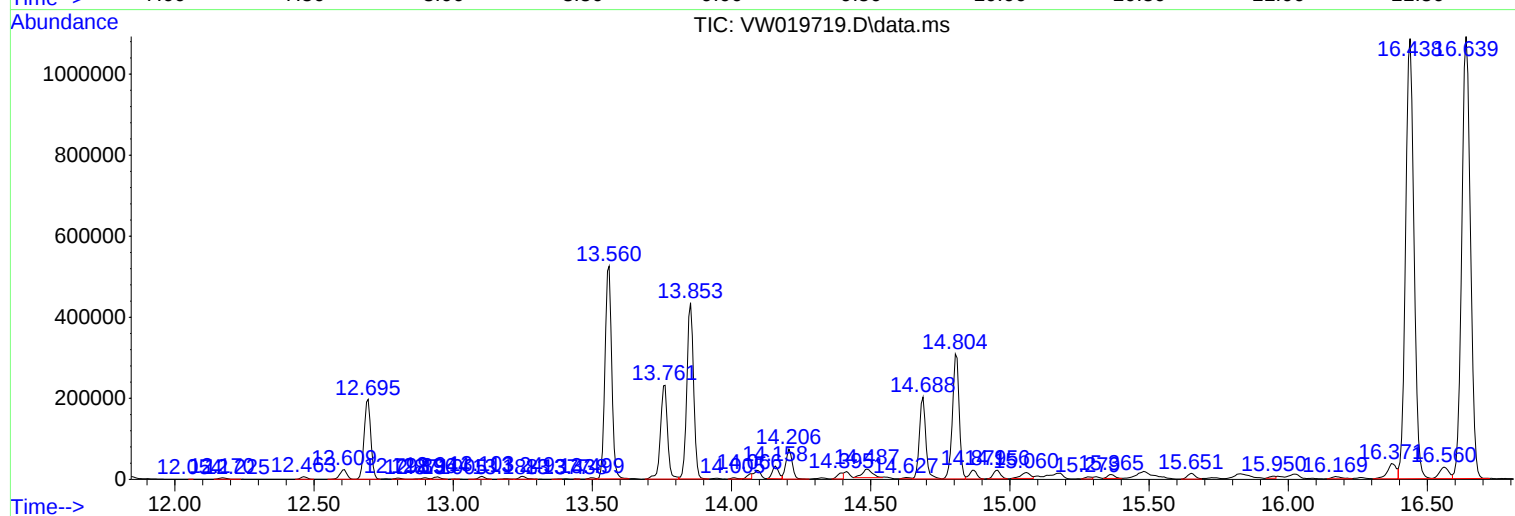
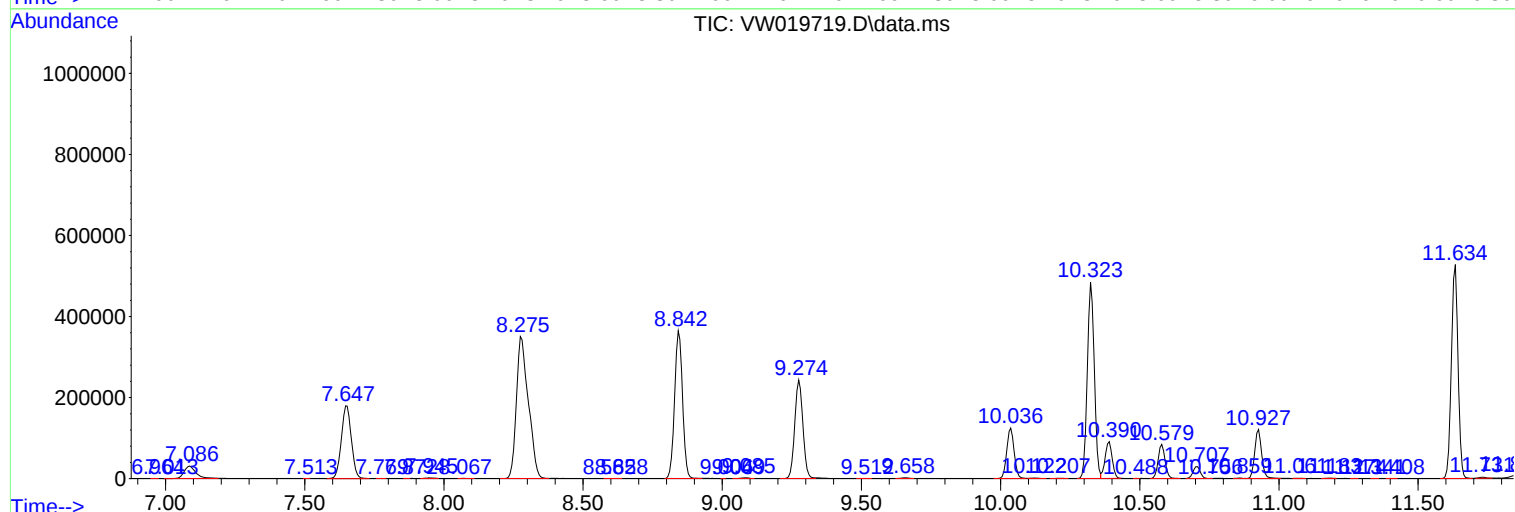
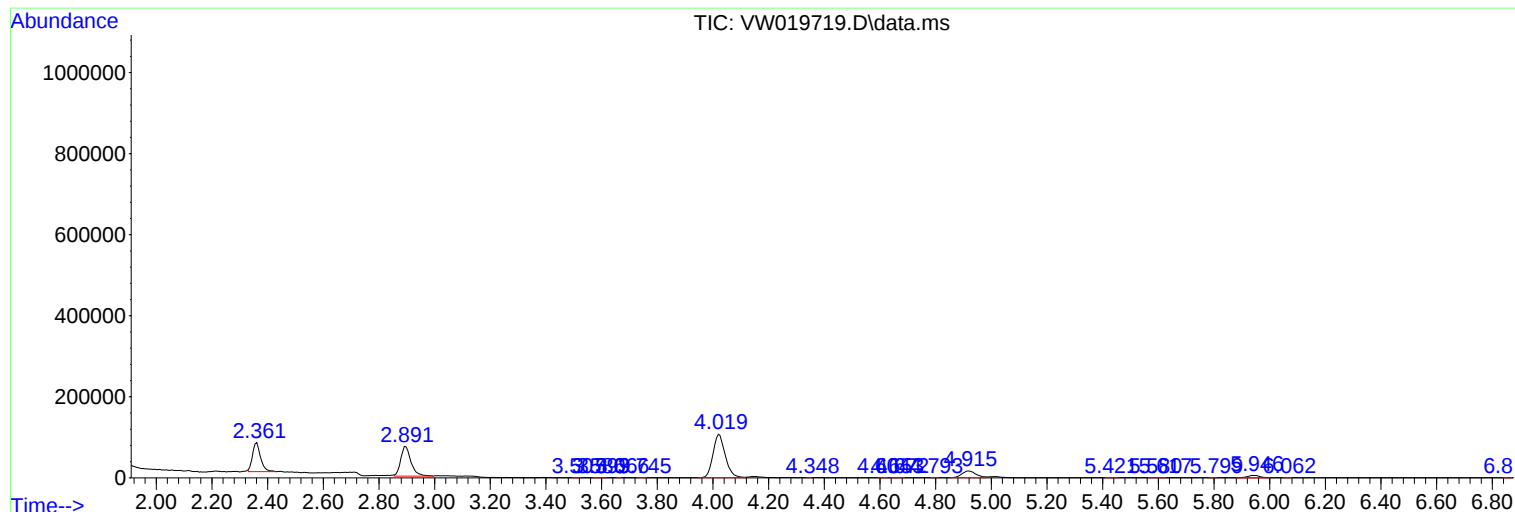
Sum of corrected areas: 14617819

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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM081121SMA.M
Quant Title : SFAM01.0

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



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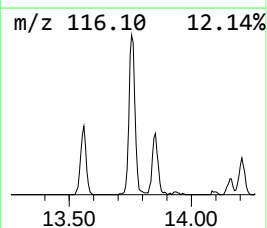
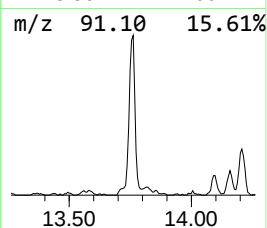
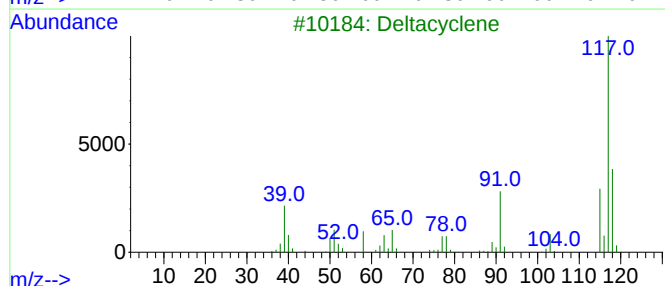
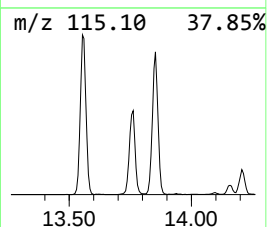
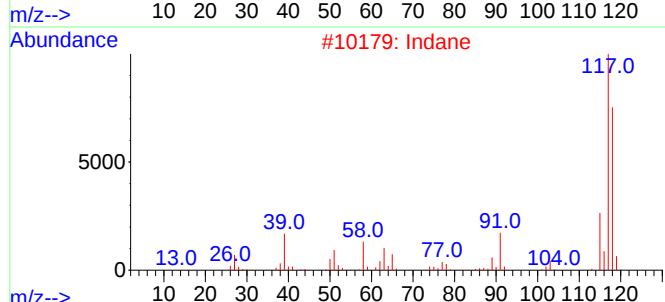
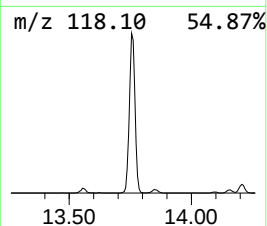
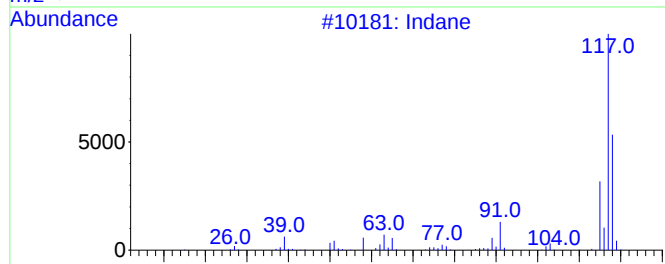
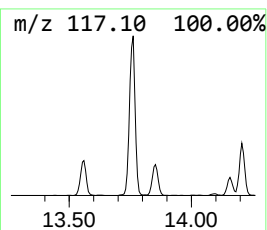
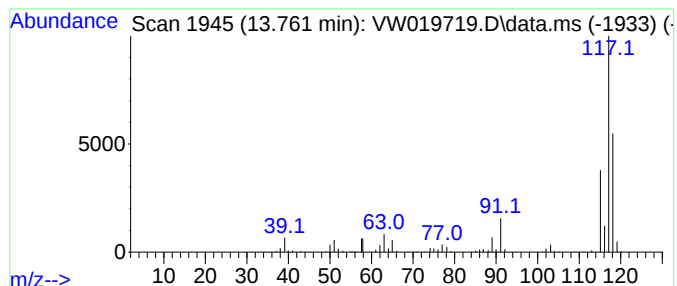
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM081121SMA.M
Quant Title : SFAM01.0

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

Peak Number 3 Indane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.761	11.70 ug/L	406477	1,4-Dichlorobenzene-d4	13.560

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane			118	C9H10	000496-11-7	93
2	Indane			118	C9H10	000496-11-7	91
3	Deltacyclene			118	C9H10	007785-10-6	80
4	Benzene, 1-ethenyl-4-methyl-			118	C9H10	000622-97-9	74
5	Benzene, cyclopropyl-			118	C9H10	000873-49-4	72



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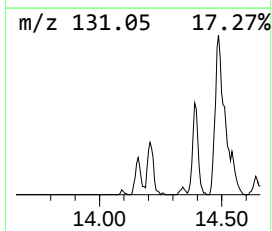
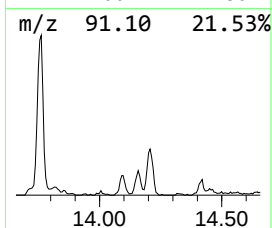
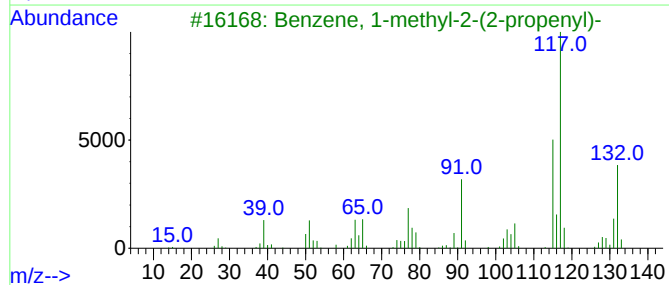
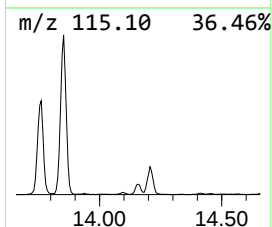
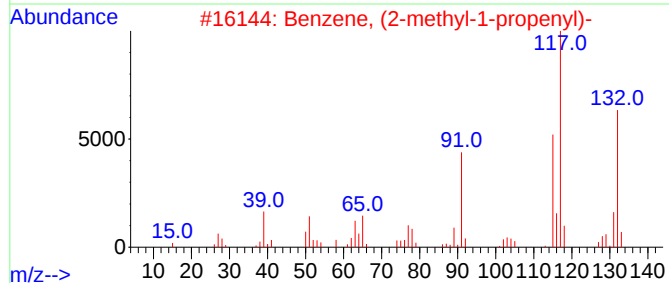
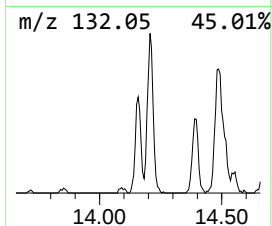
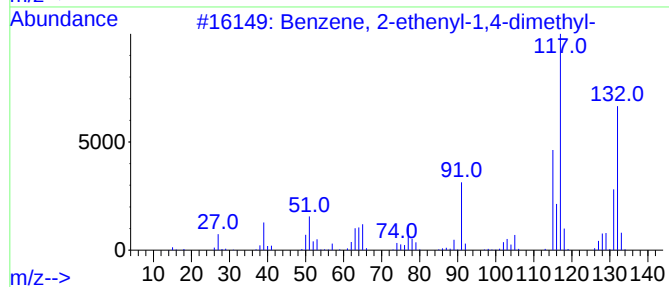
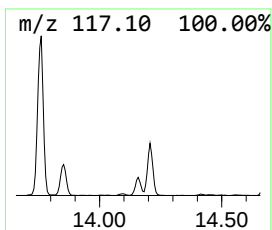
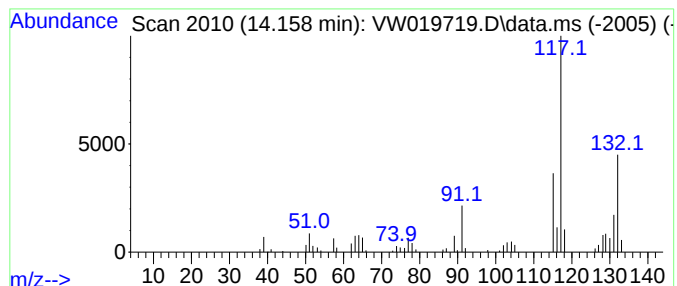
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM081121SMA.M
Quant Title : SFAM01.0

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

Peak Number 4 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.158	1.46 ug/L	50628	1,4-Dichlorobenzene-d4	13.560

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	95
2			Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	94
3			Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	94
4			Benzene, (2-methyl-2-propenyl)-	132	C10H12	003290-53-7	93
5			(E)-1-Phenyl-1-butene	132	C10H12	001005-64-7	91



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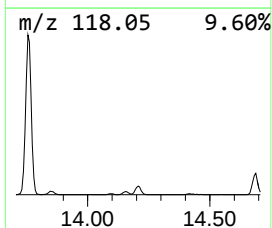
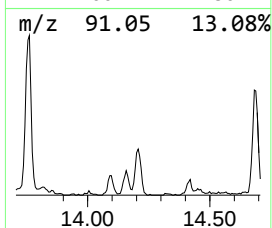
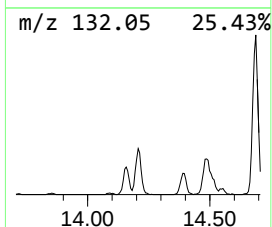
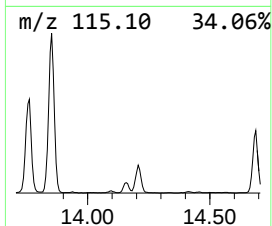
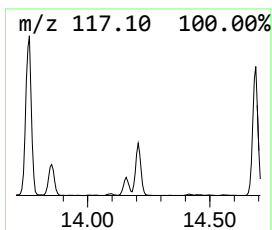
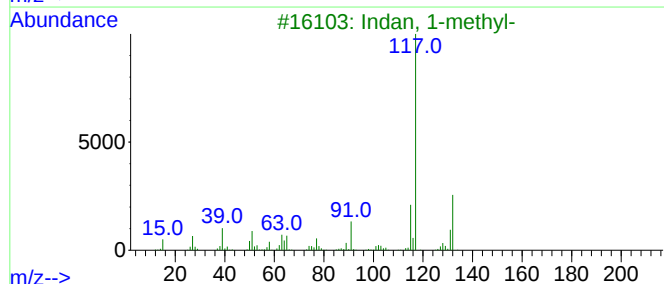
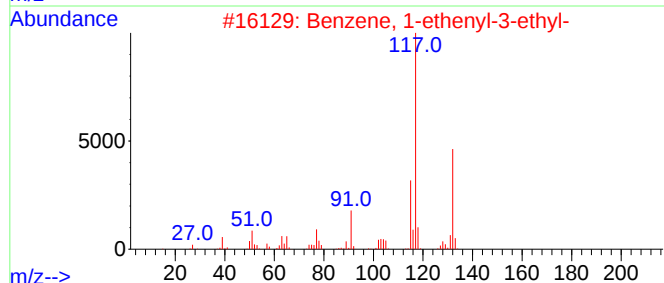
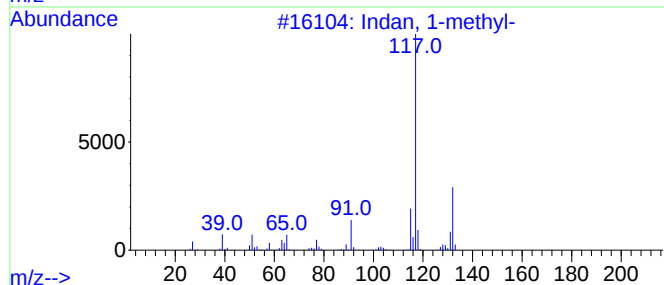
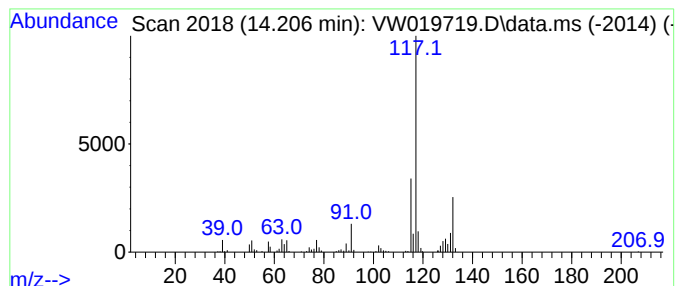
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM081121SMA.M
Quant Title : SFAM01.0

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

Peak Number 5 Indan, 1-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.206	3.32 ug/L	115219	1,4-Dichlorobenzene-d4	13.560

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indan, 1-methyl-	132	C10H12	000767-58-8	87
2			Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	86
3			Indan, 1-methyl-	132	C10H12	000767-58-8	81
4			Benzene, 2-butenyl-	132	C10H12	001560-06-1	80
5			Benzene, (2-methyl-2-propenyl)-	132	C10H12	003290-53-7	80



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW081221\
Data File : VW019719.D
Acq On : 12 Aug 2021 11:39
Operator : SY/VA
Sample : M3365-03
Misc : 5.92g/10mL/MSVOA_W/SOIL
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
DBK83

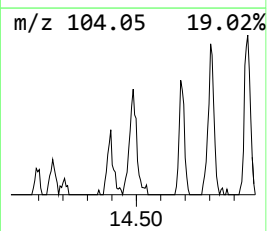
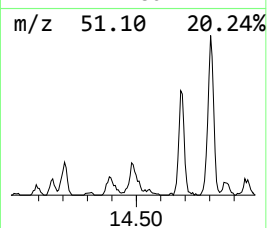
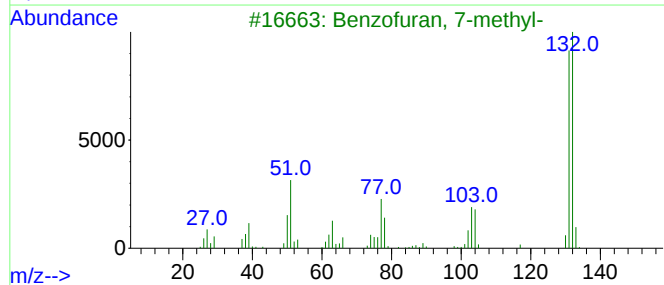
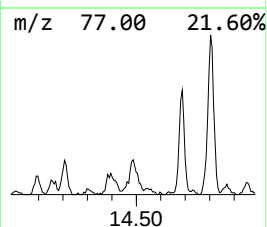
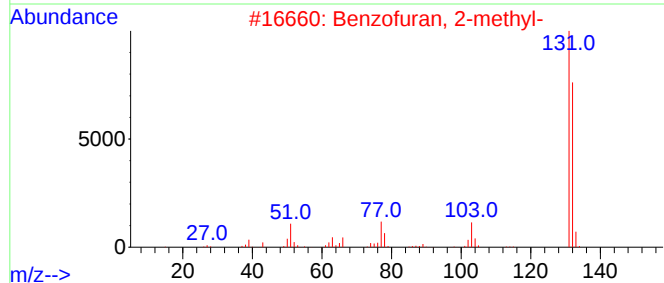
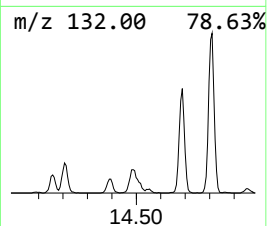
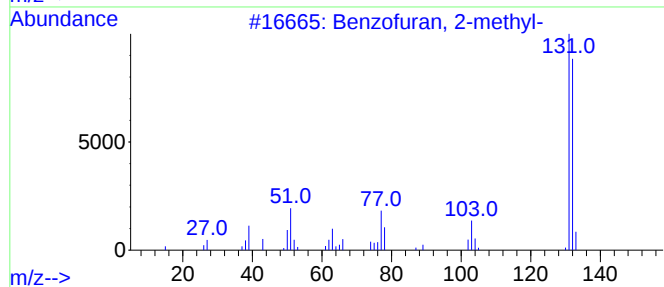
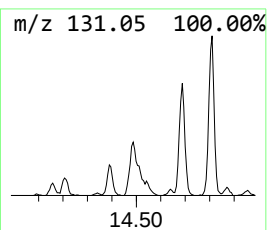
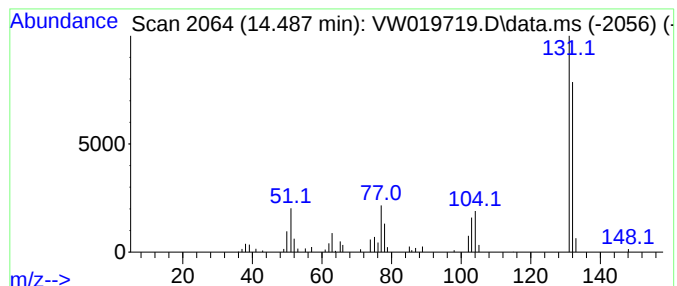
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Quant Title : SFAM01.0

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

Peak Number 6 Benzfuran, 2-methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.487	1.52 ug/L	52682	1,4-Dichlorobenzene-d4	13.560

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	94
2			Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	94
3			Benzofuran, 7-methyl-	132	C9H8O	017059-52-8	93
4			2-Propenal, 3-phenyl-	132	C9H8O	000104-55-2	91
5			Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW081221\
Data File : VW019719.D
Acq On : 12 Aug 2021 11:39
Operator : SY/VA
Sample : M3365-03
Misc : 5.92g/10mL/MSVOA_W/SOIL
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
DBK83

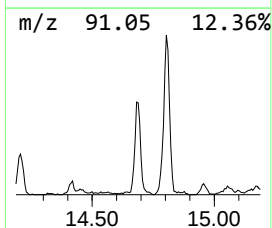
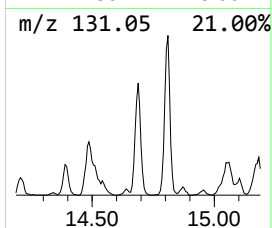
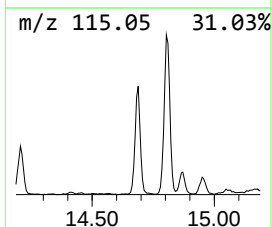
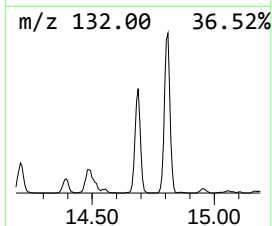
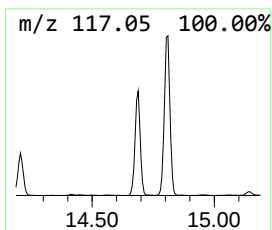
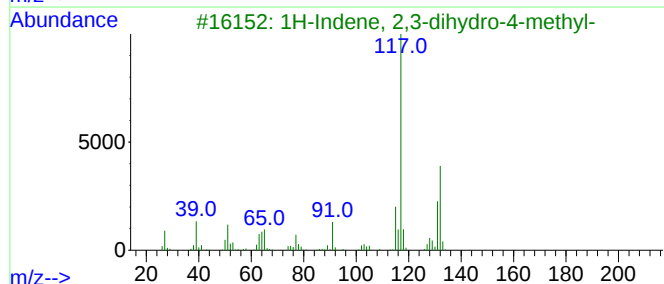
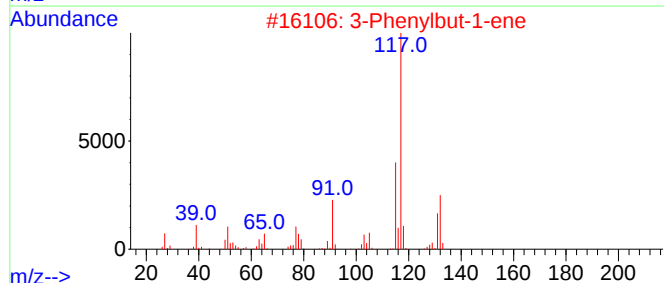
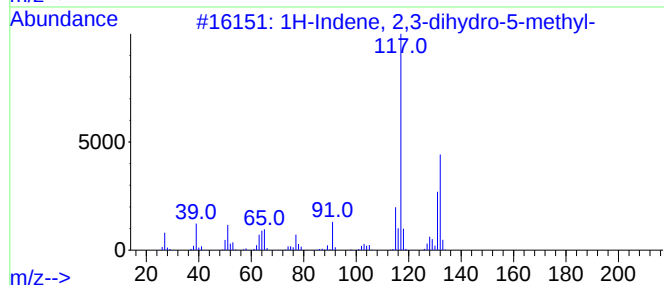
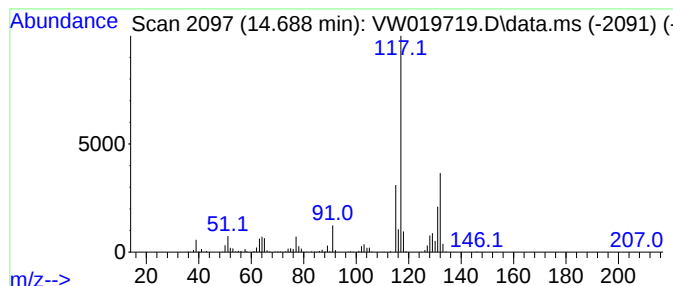
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Quant Title : SFAM01.0

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

Peak Number 7 1H-Indene, 2,3-dihydro-5-me... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.688	9.32 ug/L	323650	1,4-Dichlorobenzene-d4	13.560

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	90
2		3-Phenylbut-1-ene	132	C10H12	000934-10-1	90
3		1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	87
4		Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	87
5		Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	83



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW081221\
Data File : VW019719.D
Acq On : 12 Aug 2021 11:39
Operator : SY/VA
Sample : M3365-03
Misc : 5.92g/10mL/MSVOA_W/SOIL
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
DBK83

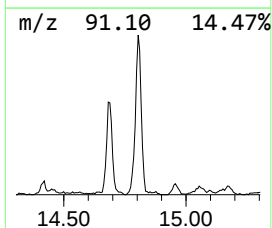
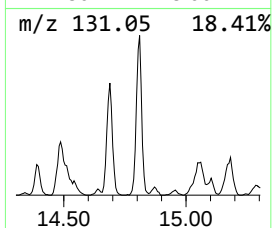
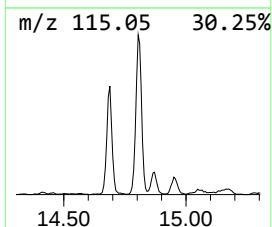
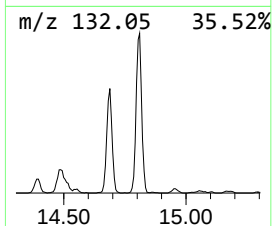
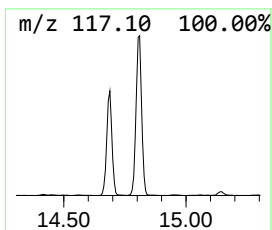
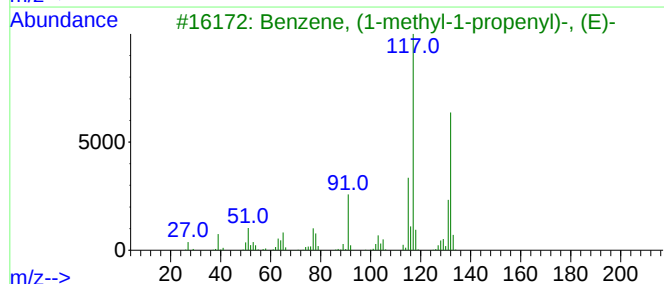
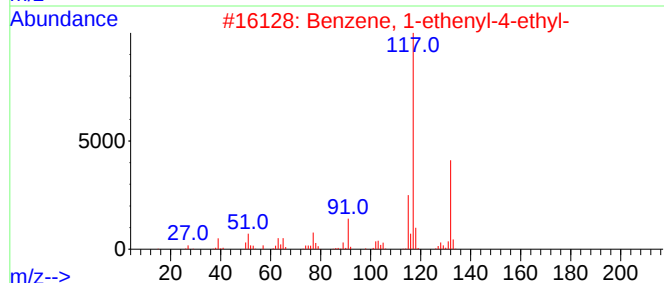
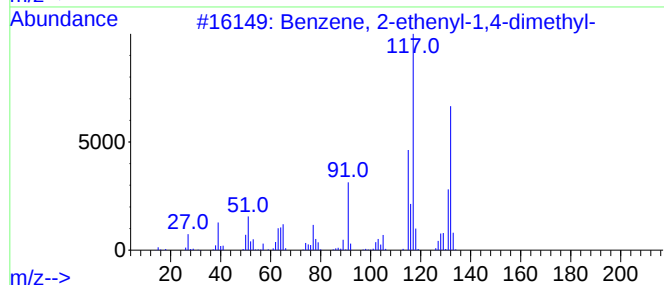
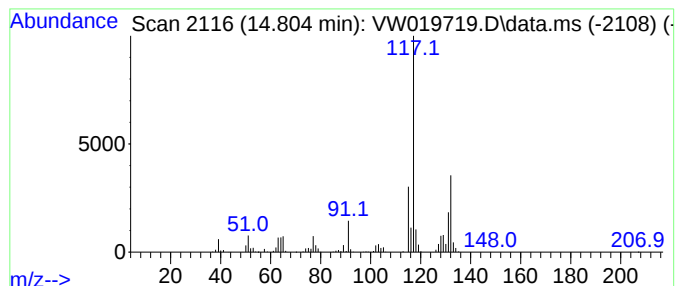
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Quant Title : SFAM01.0

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

Peak Number 8 Benzene, 1-ethenyl-4-ethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.804	15.06 ug/L	522917	1,4-Dichlorobenzene-d4	13.560

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	96
2			Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	94
3			Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	91
4			Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	90
5			1-Phenyl-1-butene	132	C10H12	000824-90-8	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW081221\
Data File : VW019719.D
Acq On : 12 Aug 2021 11:39
Operator : SY/VA
Sample : M3365-03
Misc : 5.92g/10mL/MSVOA_W/SOIL
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
DBK83

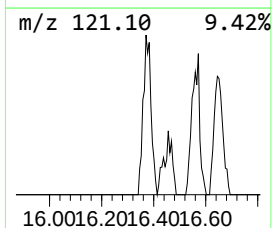
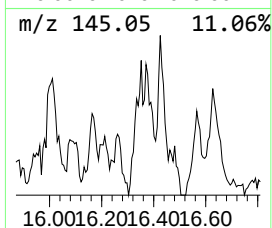
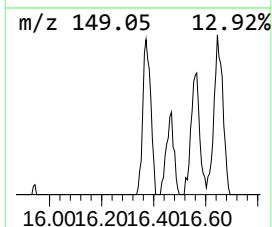
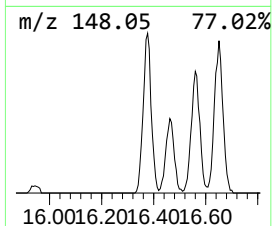
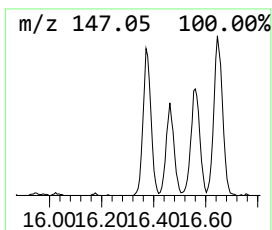
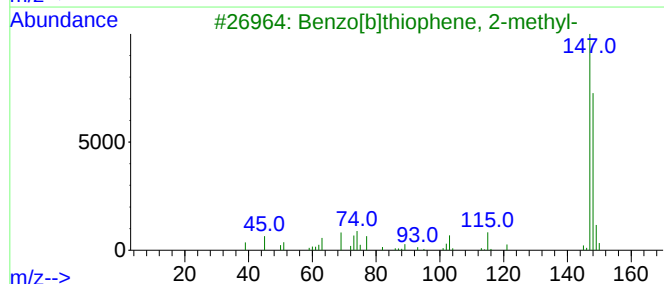
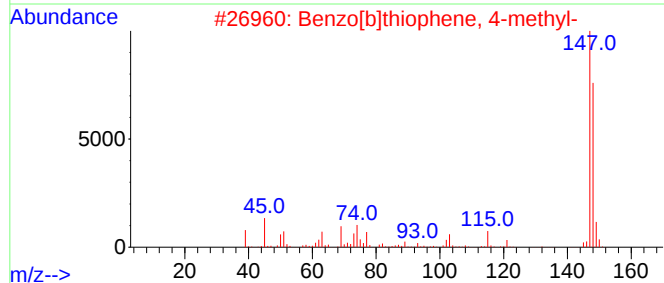
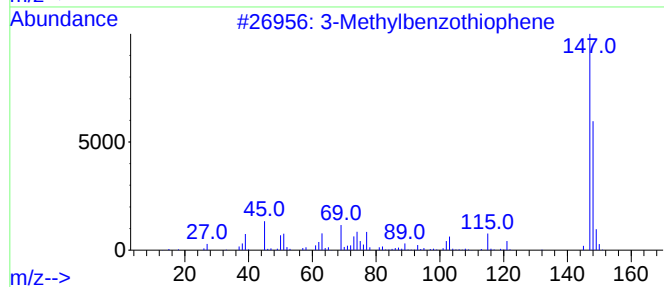
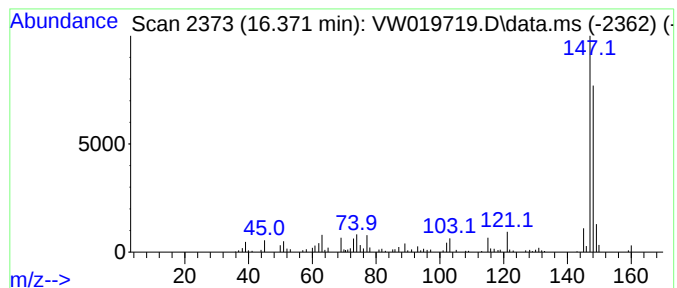
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Quant Title : SFAM01.0

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

Peak Number 9 3-Methylbenzothiophene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.371	2.60 ug/L	90345	1,4-Dichlorobenzene-d4	13.560

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Methylbenzothiophene	148	C9H8S	001455-18-1	91
2			Benzo[b]thiophene, 4-methyl-	148	C9H8S	014315-11-8	90
3			Benzo[b]thiophene, 2-methyl-	148	C9H8S	001195-14-8	90
4			Benzo[b]thiophene, 2-methyl-	148	C9H8S	001195-14-8	90
5			Benzo[b]thiophene, 5-methyl-	148	C9H8S	014315-14-1	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW081221\
Data File : VW019719.D
Acq On : 12 Aug 2021 11:39
Operator : SY/VA
Sample : M3365-03
Misc : 5.92g/10mL/MSVOA_W/SOIL
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
DBK83

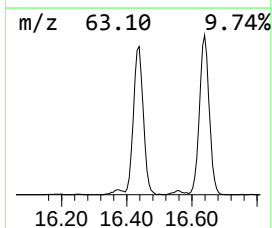
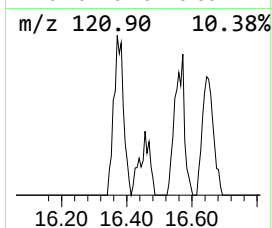
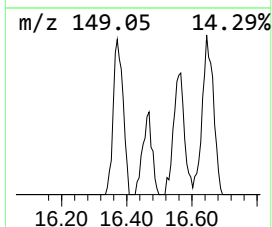
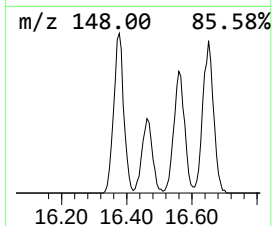
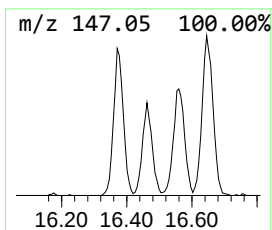
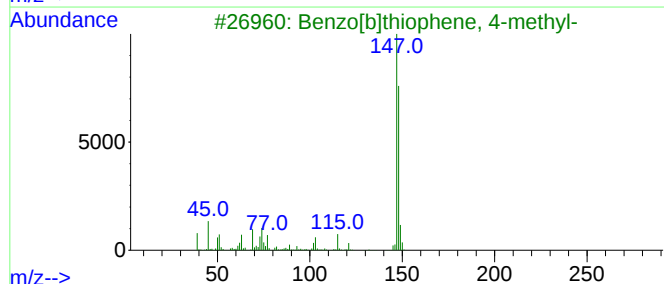
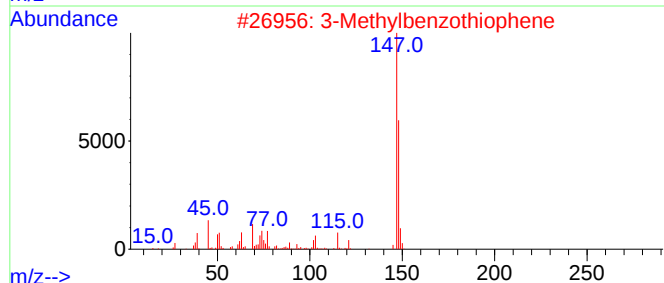
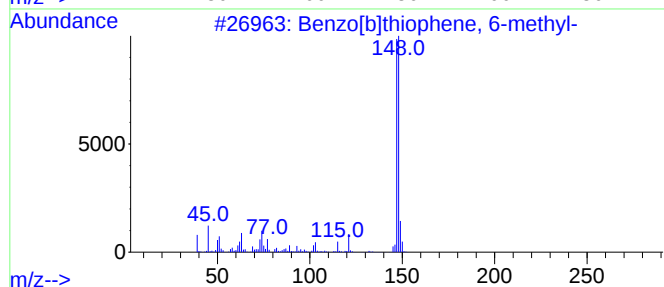
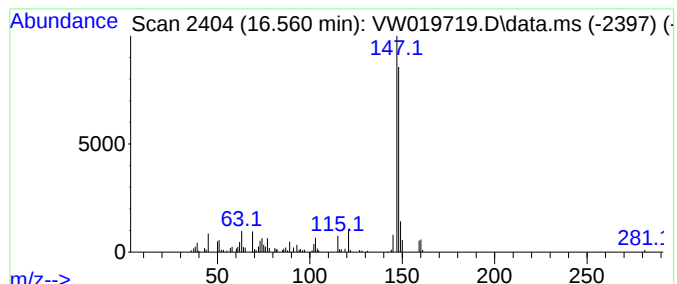
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Quant Title : SFAM01.0

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

Peak Number 11 Benzo[b]thiophene, 6-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.560	1.95 ug/L	67708	1,4-Dichlorobenzene-d4	13.560

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[b]thiophene, 6-methyl-	148	C9H8S	016587-47-6	91
2			3-Methylbenzothiophene	148	C9H8S	001455-18-1	90
3			Benzo[b]thiophene, 4-methyl-	148	C9H8S	014315-11-8	87
4			Benzo[b]thiophene, 2-methyl-	148	C9H8S	001195-14-8	87
5			Benzo[b]thiophene, 5-methyl-	148	C9H8S	014315-14-1	87



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW081221\
Data File : VW019719.D
Acq On : 12 Aug 2021 11:39
Operator : SY/VA
Sample : M3365-03
Misc : 5.92g/10mL/MSVOA_W/SOIL
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
DBK83

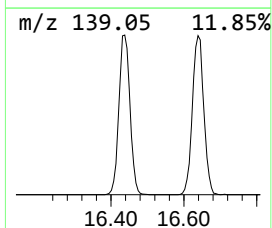
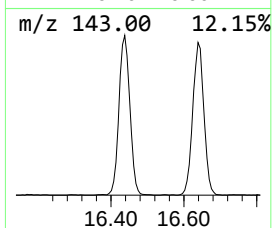
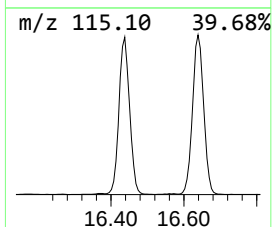
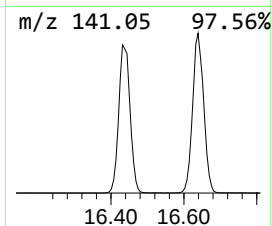
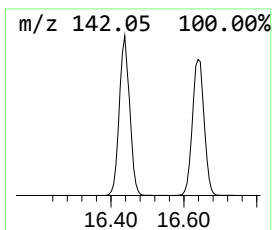
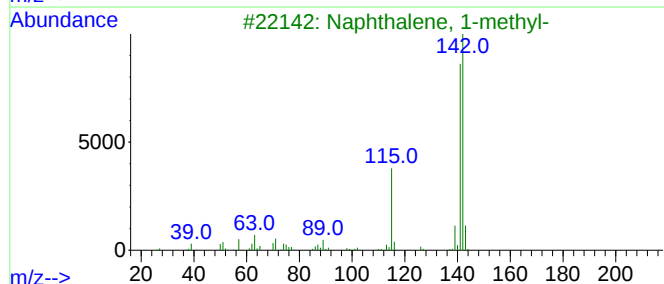
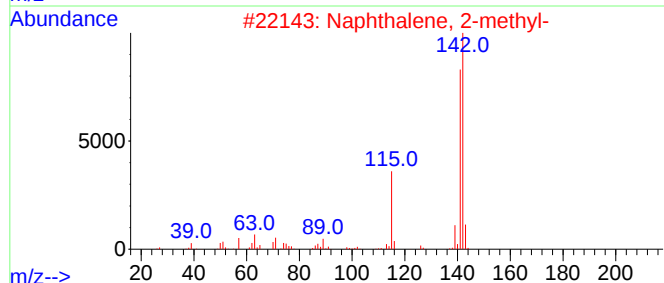
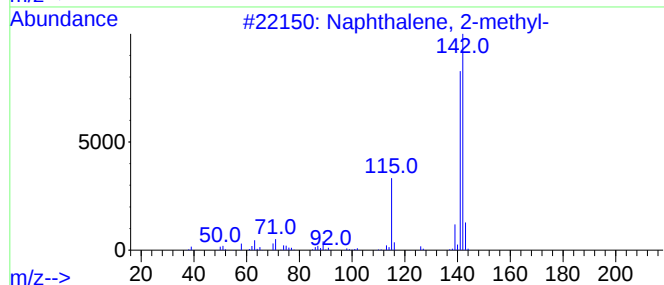
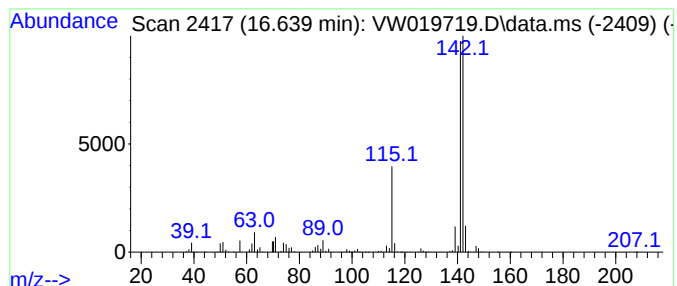
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Quant Title : SFAM01.0

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

Peak Number 12 1,4-Methanonaphthalene, 1,4... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.639	68.20 ug/L	2368580	1,4-Dichlorobenzene-d4	13.560

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
2			Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
3			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
4			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	95
5			1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW081221\
Data File : VW019719.D
Acq On : 12 Aug 2021 11:39
Operator : SY/VA
Sample : M3365-03
Misc : 5.92g/10mL/MSVOA_W/SOIL
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
DBK83

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM081121SMA.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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Benzene, 2-ethe...	14.158	1.5	ug/L	50628	3	13.560	868267	25.0
Indan, 1-methyl-	14.206	3.3	ug/L	115219	3	13.560	868267	25.0
Benzofuran, 2-m...	14.487	1.5	ug/L	52682	3	13.560	868267	25.0
1H-Indene, 2,3-...	14.688	9.3	ug/L	323650	3	13.560	868267	25.0
Benzene, 1-ethe...	14.804	15.1	ug/L	522917	3	13.560	868267	25.0
3-Methylbenzoth...	16.371	2.6	ug/L	90345	3	13.560	868267	25.0
Benzo[b]thiophe...	16.560	2.0	ug/L	67708	3	13.560	868267	25.0
1,4-Methanonaph...	16.639	68.2	ug/L	2368580	3	13.560	868267	25.0