

Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW120221\
 Data File : VW021052.D
 Acq On : 02 Dec 2021 17:13
 Operator : SY/VA
 Sample : M4886-18
 Misc : 5.00g/10.0mL/MSVOA_W/SOIL
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 EX897

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\SFAMWLM111521SMA.M
 Title : SFAM01.0

Signal : TIC: VW021052.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.166	41	43	47	rVB2	815	683	0.21%	0.023%
2	2.203	47	49	50	rBV	543	401	0.12%	0.013%
3	2.349	67	73	81	rVB	25869	45107	13.82%	1.492%
4	2.824	149	151	153	rVB2	522	552	0.17%	0.018%
5	2.885	153	161	172	rBV2	14773	32822	10.06%	1.086%
6	3.379	234	242	247	rVB3	469	1166	0.36%	0.039%
7	4.013	336	346	359	rBV3	26007	80787	24.76%	2.673%
8	4.147	364	368	373	rBV2	692	1520	0.47%	0.050%
9	4.379	396	406	414	rBV2	1515	4539	1.39%	0.150%
10	4.641	447	449	453	rBV	92	173	0.05%	0.006%
11	4.781	464	472	481	rBV3	1232	3653	1.12%	0.121%
12	4.909	484	493	494	rBV2	4107	7085	2.17%	0.234%
13	5.092	519	523	525	rBV	156	216	0.07%	0.007%
14	5.415	573	576	579	rBV3	320	474	0.15%	0.016%
15	5.641	611	613	619	rBB	162	296	0.09%	0.010%
16	5.720	619	626	627	rBV	150	282	0.09%	0.009%
17	5.787	633	637	638	rBV2	154	192	0.06%	0.006%
18	5.934	653	661	670	rBB3	874	2410	0.74%	0.080%
19	6.013	671	674	677	rBV	112	186	0.06%	0.006%
20	6.159	696	698	701	rBV	154	155	0.05%	0.005%
21	6.348	725	729	733	rVV2	117	201	0.06%	0.007%
22	6.970	826	831	835	rBB2	191	333	0.10%	0.011%
23	7.086	840	850	856	rBV	8236	22901	7.02%	0.758%
24	7.159	858	862	874	rVB4	5664	15560	4.77%	0.515%
25	7.403	901	902	906	rVB	146	166	0.05%	0.005%
26	7.647	932	942	956	rBV	53321	130831	40.10%	4.329%
27	7.903	981	984	986	rBV	130	170	0.05%	0.006%
28	8.025	1001	1004	1008	rBV	261	359	0.11%	0.012%
29	8.275	1030	1045	1061	rBV2	101420	312292	95.71%	10.333%
30	8.409	1061	1067	1072	rBV3	1191	2305	0.71%	0.076%
31	8.470	1075	1077	1079	rBV	287	194	0.06%	0.006%
32	8.531	1085	1087	1090	rVB2	200	211	0.06%	0.007%
33	8.842	1128	1138	1150	rVB	134953	270910	83.03%	8.963%
34	9.092	1165	1179	1185	rBV4	3801	7853	2.41%	0.260%
35	9.195	1191	1196	1197	rBV	127	206	0.06%	0.007%
36	9.220	1197	1200	1201	rBV2	341	279	0.09%	0.009%

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Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM111521SMA.M
 Title : SFAM01.0

37	9.275	1201	1209	1217	rBV	71116	142388	43.64%	4.711%
38	9.500	1245	1246	1251	rVB	152	175	0.05%	0.006%
39	9.610	1262	1264	1267	rBV	206	232	0.07%	0.008%
40	9.659	1270	1272	1277	rVB	398	526	0.16%	0.017%
41	9.750	1284	1287	1290	rBV	464	541	0.17%	0.018%
42	9.890	1305	1310	1315	rVB4	383	730	0.22%	0.024%
43	9.982	1321	1325	1326	rBV	222	177	0.05%	0.006%
44	10.037	1326	1334	1345	rVB	33807	61974	18.99%	2.050%
45	10.128	1347	1349	1351	rVB	280	186	0.06%	0.006%
46	10.146	1351	1352	1358	rVB	159	211	0.06%	0.007%
47	10.213	1358	1363	1366	rBV2	771	1323	0.41%	0.044%
48	10.323	1373	1381	1388	rBV	142710	253037	77.55%	8.372%
49	10.390	1388	1392	1398	rVB2	3306	5768	1.77%	0.191%
50	10.439	1398	1400	1401	rVB	233	161	0.05%	0.005%
51	10.494	1402	1409	1416	rVB3	577	1259	0.39%	0.042%
52	10.579	1416	1423	1430	rBV	21169	36938	11.32%	1.222%
53	10.665	1433	1437	1438	rBV2	470	563	0.17%	0.019%
54	10.701	1438	1443	1448	rVB2	4996	8089	2.48%	0.268%
55	10.762	1451	1453	1458	rVB2	629	873	0.27%	0.029%
56	10.860	1464	1469	1474	rBV2	3142	5634	1.73%	0.186%
57	10.921	1474	1479	1490	rBV	36643	62553	19.17%	2.070%
58	11.024	1490	1496	1497	rBV4	546	933	0.29%	0.031%
59	11.067	1497	1503	1511	rBV2	5146	9520	2.92%	0.315%
60	11.128	1511	1513	1514	rBV	438	328	0.10%	0.011%
61	11.177	1519	1521	1526	rVB4	577	636	0.19%	0.021%
62	11.225	1526	1529	1534	rBV6	975	1773	0.54%	0.059%
63	11.268	1534	1536	1537	rBV2	487	422	0.13%	0.014%
64	11.286	1537	1539	1540	rBV2	460	356	0.11%	0.012%
65	11.347	1543	1549	1550	rBV3	1290	2014	0.62%	0.067%
66	11.433	1559	1563	1569	rVB2	5782	10552	3.23%	0.349%
67	11.628	1587	1595	1603	rBV	197042	326293	100.00%	10.796%
68	11.725	1607	1611	1616	rVB2	2917	4619	1.42%	0.153%
69	11.835	1624	1629	1635	rBV	2607	4602	1.41%	0.152%
70	11.945	1643	1647	1648	rBV2	334	264	0.08%	0.009%
71	11.963	1648	1650	1652	rBV3	485	381	0.12%	0.013%
72	12.030	1659	1661	1667	rVB4	729	907	0.28%	0.030%
73	12.170	1680	1684	1690	rVB7	1695	3330	1.02%	0.110%
74	12.554	1745	1747	1749	rVV3	693	540	0.17%	0.018%
75	12.603	1749	1755	1763	rVV2	15581	29182	8.94%	0.966%
76	12.689	1763	1769	1776	rVB	61223	103114	31.60%	3.412%

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77	12.878	1787	1800	1801	rBV2	13529	39984	12.25%	1.323%
78	13.292	1856	1868	1869	rBV5	21669	44764	13.72%	1.481%
79	13.554	1905	1911	1919	rVB	202845	324033	99.31%	10.721%
80	13.652	1922	1927	1932	rBV2	5330	9355	2.87%	0.310%
81	13.853	1953	1960	1969	rBV	148103	255892	78.42%	8.467%
82	14.066	1990	1995	1999	rBV2	7951	13972	4.28%	0.462%
83	14.200	2015	2017	2022	rBV3	1733	2310	0.71%	0.076%
84	14.389	2045	2048	2049	rBV3	2234	2201	0.67%	0.073%
85	14.950	2130	2140	2152	rVB2	45712	143181	43.88%	4.737%
86	15.066	2153	2159	2168	rVV5	7567	17083	5.24%	0.565%
87	15.280	2190	2194	2197	rBV2	5761	9121	2.80%	0.302%
88	15.359	2203	2207	2213	rVB	14572	27196	8.33%	0.900%
89	15.432	2216	2219	2223	rBV4	5625	9059	2.78%	0.300%
90	15.633	2243	2252	2264	rVB4	9052	33103	10.15%	1.095%
91	15.779	2274	2276	2277	rBV	624	412	0.13%	0.014%
92	15.840	2283	2286	2291	rVB5	2650	3911	1.20%	0.129%
93	15.895	2291	2295	2296	rBV3	1904	2706	0.83%	0.090%
94	16.023	2314	2316	2322	rVB2	2448	3115	0.95%	0.103%
95	16.096	2325	2328	2329	rBV3	1911	2162	0.66%	0.072%
96	16.261	2352	2355	2362	rVB5	2308	4072	1.25%	0.135%
97	16.371	2362	2373	2377	rBV7	3638	12614	3.87%	0.417%
98	16.529	2396	2399	2400	rBV2	401	372	0.11%	0.012%
99	16.639	2410	2417	2427	rVB3	13801	30488	9.34%	1.009%
100	16.724	2428	2431	2433	rBV5	772	721	0.22%	0.024%

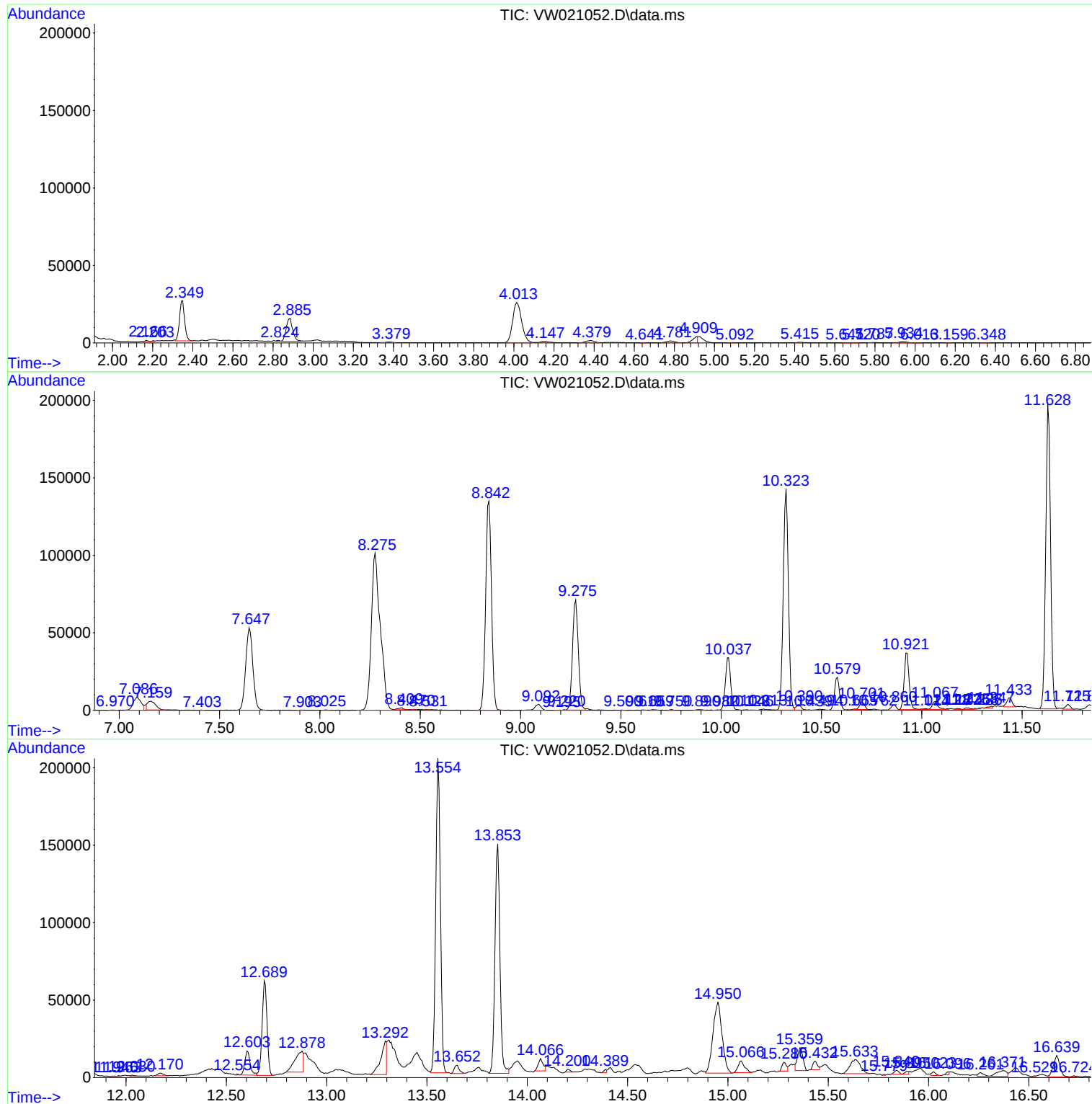
Sum of corrected areas: 3022401

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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM111521SMA.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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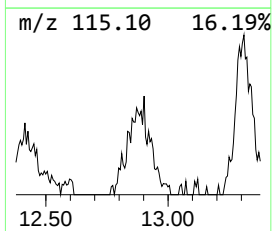
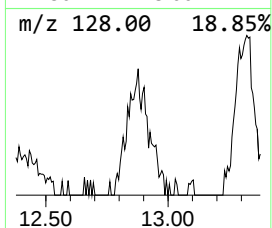
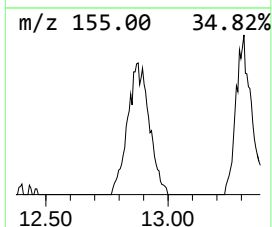
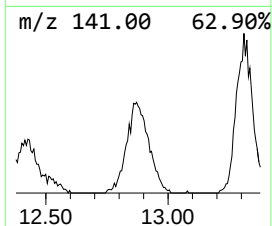
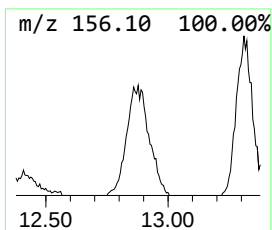
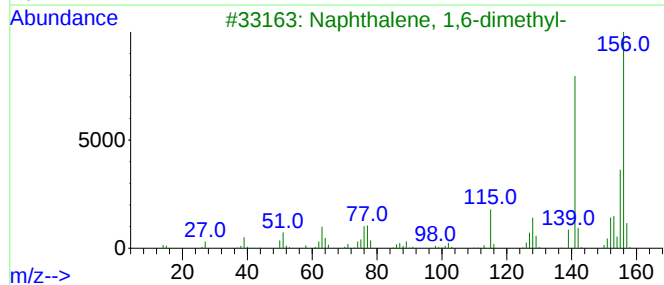
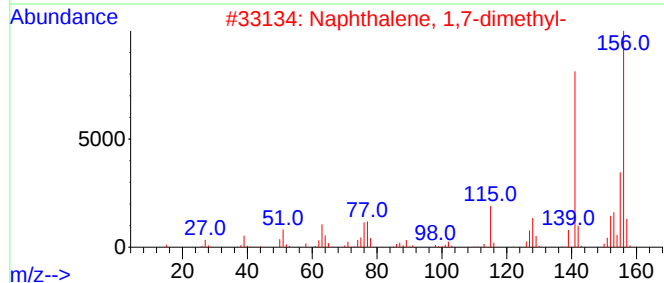
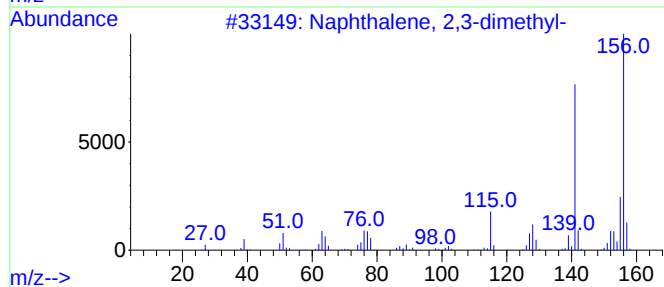
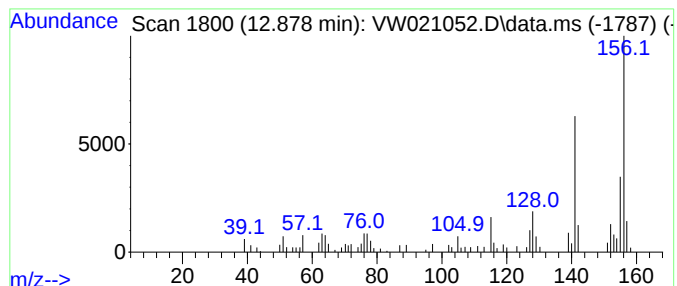
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM111521SMA.M
Quant Title : SFAM01.0

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

Peak Number 3 Naphthalene, 2,3-dimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.878	3.08 ug/L	39984	1,4-Dichlorobenzene-d4	13.554

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
2			Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	95
3			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	95
4			Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	95
5			Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	95



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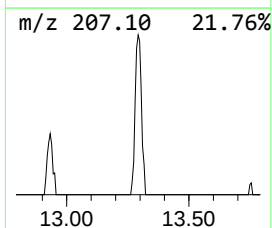
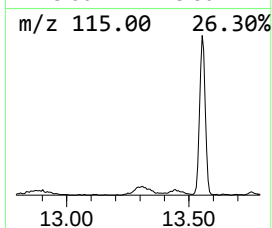
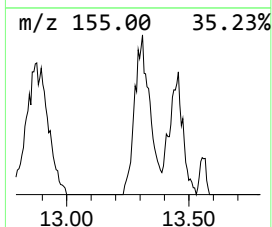
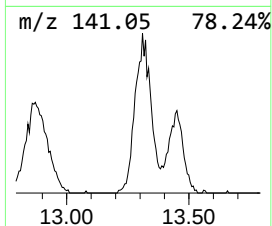
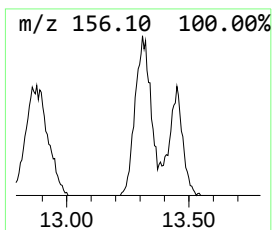
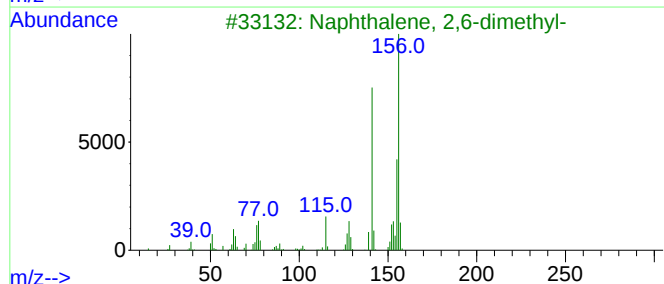
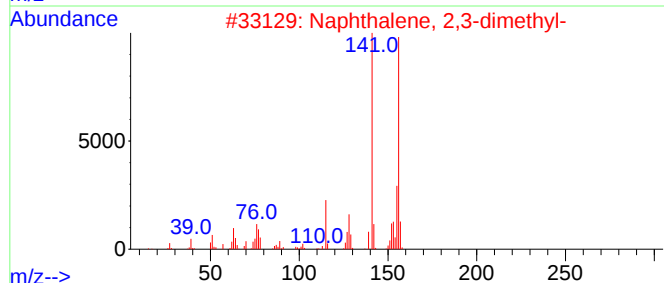
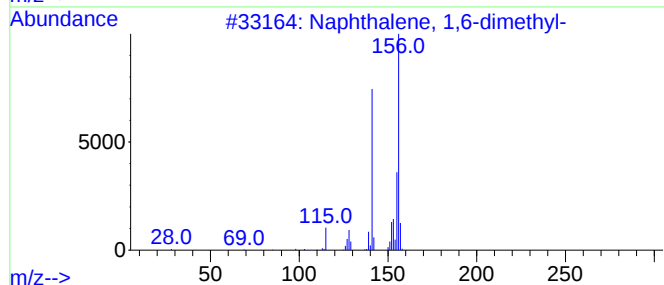
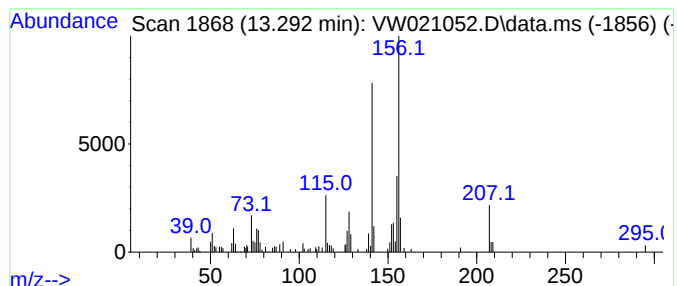
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM111521SMA.M
Quant Title : SFAM01.0

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

Peak Number 4 Naphthalene, 1,6-dimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.292	3.45 ug/L	44764	1,4-Dichlorobenzene-d4	13.554

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	96
2			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	96
3			Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	95
4			Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	95
5			Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	95



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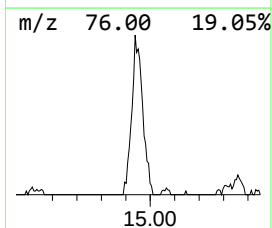
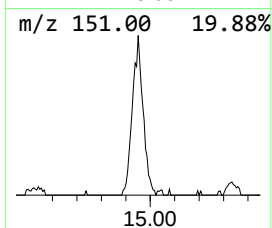
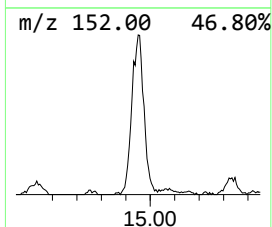
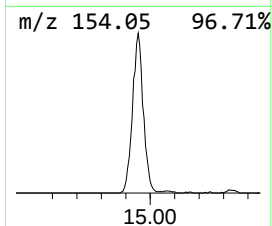
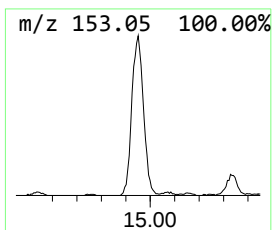
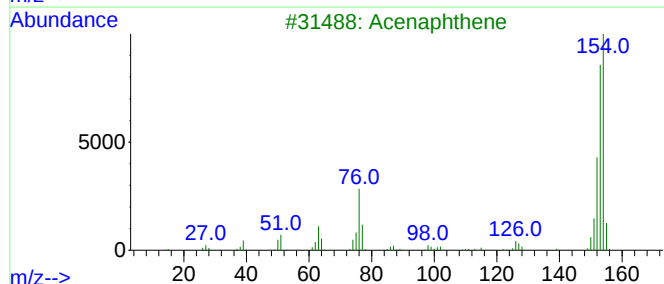
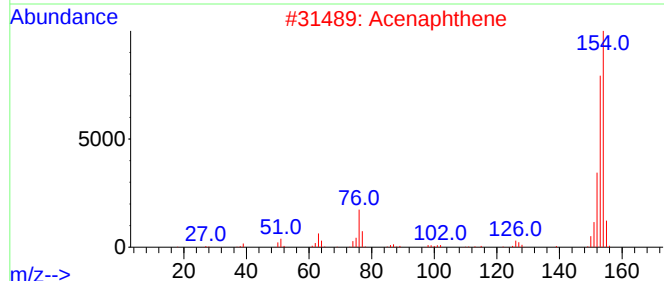
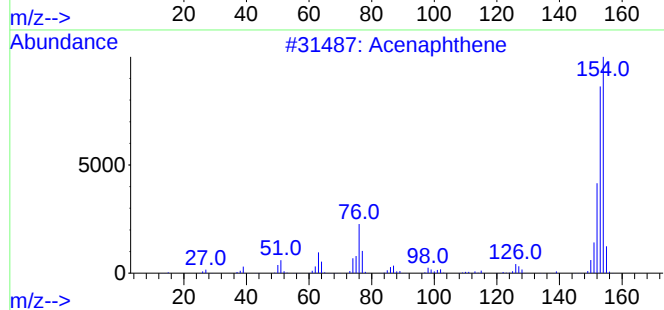
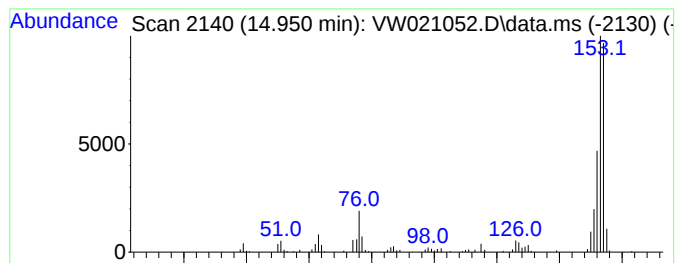
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM111521SMA.M
Quant Title : SFAM01.0

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

Peak Number 5 Acenaphthene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.950	11.05 ug/L	143181	1,4-Dichlorobenzene-d4	13.554

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Acenaphthene	154	C12H10	000083-32-9	90
2			Acenaphthene	154	C12H10	000083-32-9	81
3			Acenaphthene	154	C12H10	000083-32-9	81
4			Acenaphthene	154	C12H10	000083-32-9	70
5			Naphthalene, 2-ethenyl-	154	C12H10	000827-54-3	68



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW120221\
Data File : VW021052.D
Acq On : 02 Dec 2021 17:13
Operator : SY/VA
Sample : M4886-18
Misc : 5.00g/10.0mL/MSVOA_W/SOIL
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
EX897

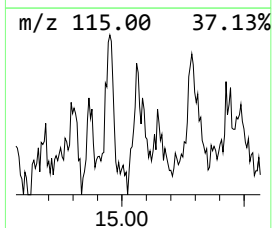
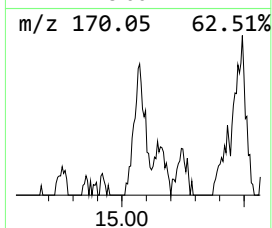
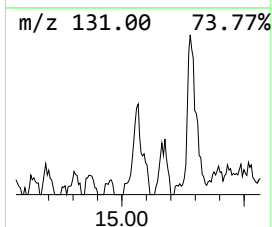
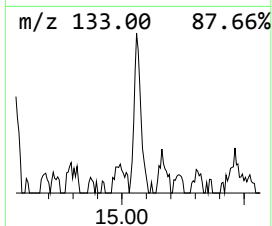
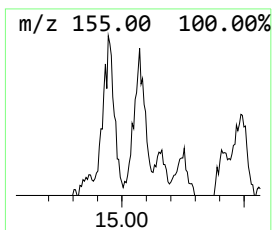
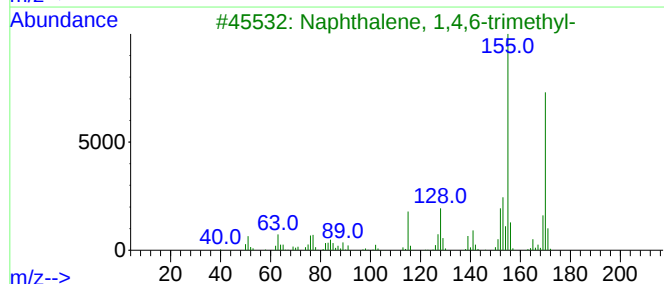
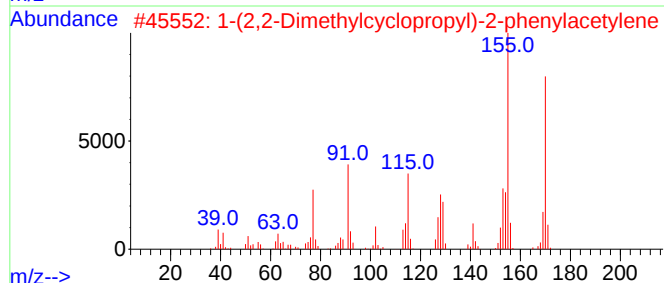
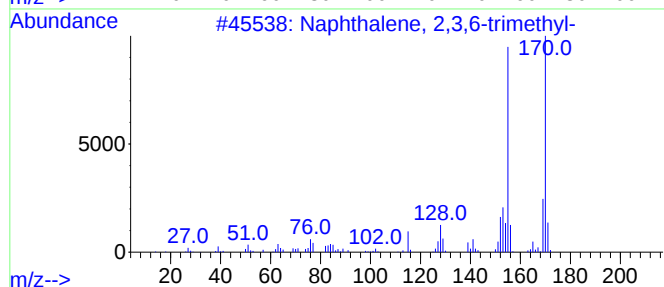
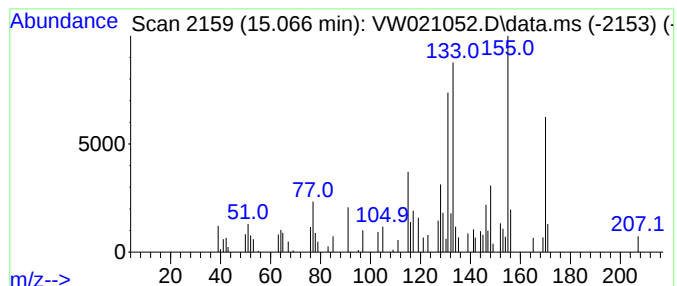
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Quant Title : SFAM01.0

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

Peak Number 6 unknown-01 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.066	1.32 ug/L	17083	1,4-Dichlorobenzene-d4	13.554

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	42
2			1-(2,2-Dimethylcyclopropyl)-2-ph...	170	C13H14	1000294-91-1	42
3			Naphthalene, 1,4,6-trimethyl-	170	C13H14	002131-42-2	38
4			3-(2-Methyl-propenyl)-1H-indene	170	C13H14	1000187-78-5	30
5			5-Methyl-1-phenylhexa-1,3,4-triene	170	C13H14	103240-79-5	25



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW120221\
Data File : VW021052.D
Acq On : 02 Dec 2021 17:13
Operator : SY/VA
Sample : M4886-18
Misc : 5.00g/10.0mL/MSVOA_W/SOIL
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
EX897

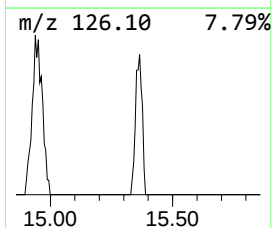
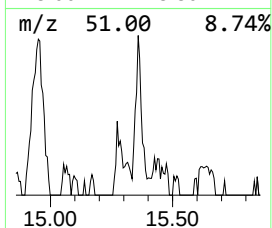
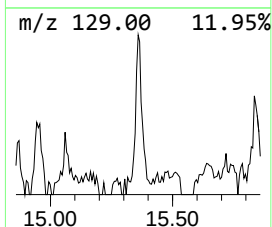
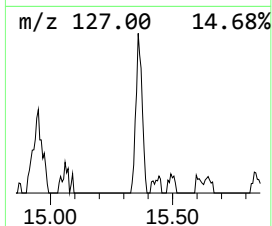
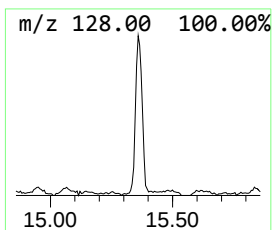
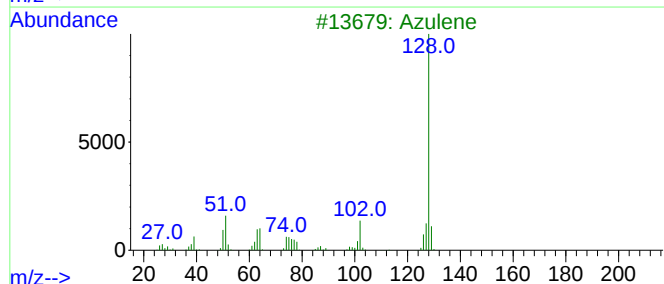
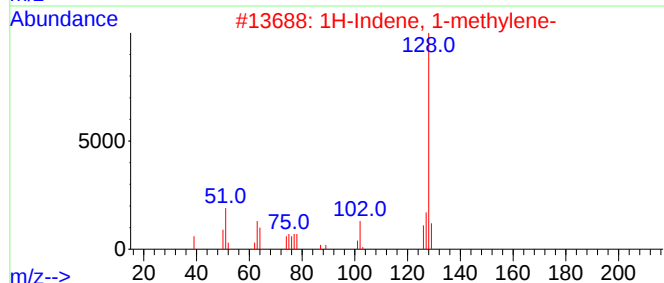
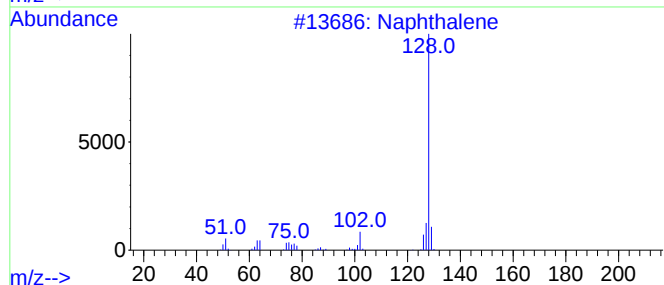
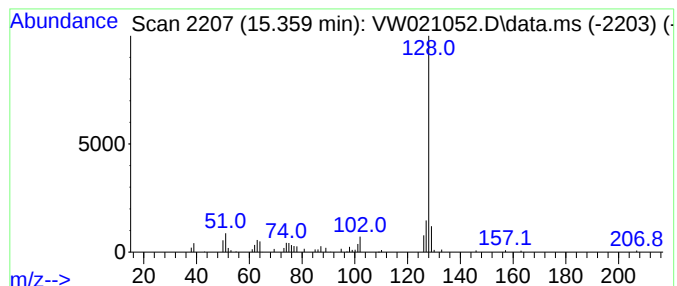
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Quant Title : SFAM01.0

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

Peak Number 7 Azulene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.359	2.10 ug/L	27196	1,4-Dichlorobenzene-d4	13.554

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene	128	C10H8	000091-20-3	93
2			1H-Indene, 1-methylene-	128	C10H8	002471-84-3	91
3			Azulene	128	C10H8	000275-51-4	91
4			Naphthalene	128	C10H8	000091-20-3	90
5			Azulene	128	C10H8	000275-51-4	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW120221\
Data File : VW021052.D
Acq On : 02 Dec 2021 17:13
Operator : SY/VA
Sample : M4886-18
Misc : 5.00g/10.0mL/MSVOA_W/SOIL
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
EX897

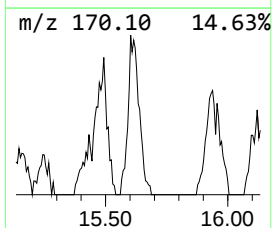
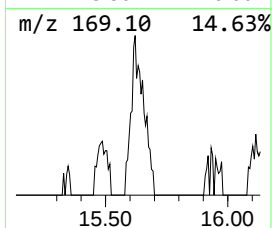
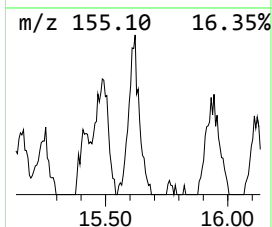
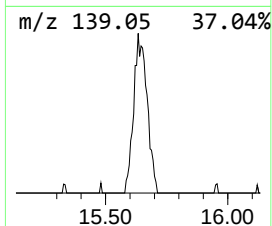
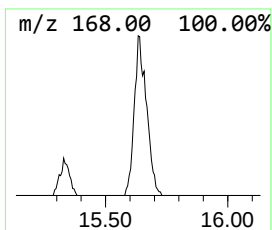
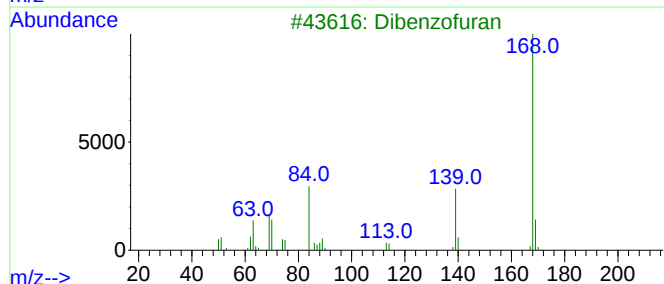
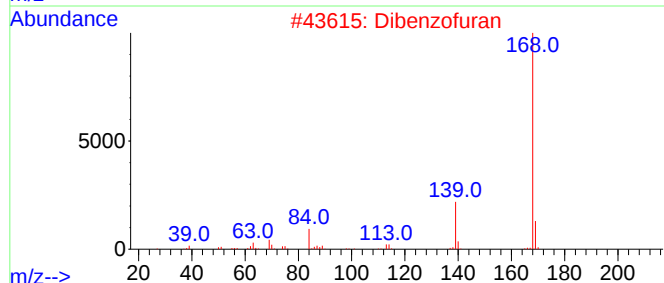
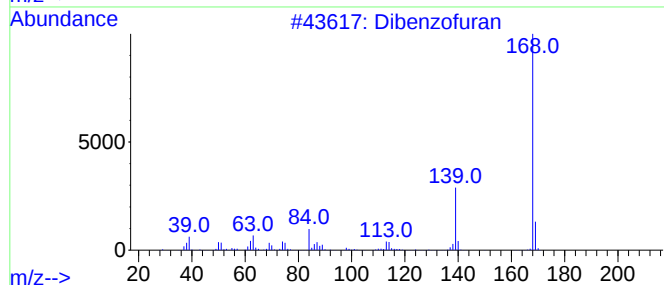
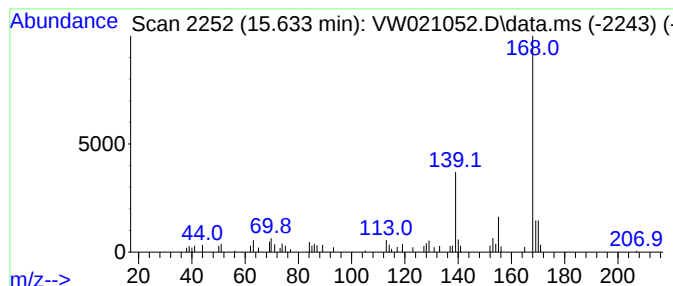
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Quant Title : SFAM01.0

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

Peak Number 8 Dibenzofuran Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.633	2.55 ug/L	33103	1,4-Dichlorobenzene-d4	13.554

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dibenzofuran	168	C12H8O	000132-64-9	68
2			Dibenzofuran	168	C12H8O	000132-64-9	64
3			Dibenzofuran	168	C12H8O	000132-64-9	59
4			3H-[1,2,4]thiadiazolo[4,3-a]pyri...	168	C6H4N2S2	1010404-55-6	52
5			Benzofuran, 5-chloro-2,3-dihydro...	168	C9H9ClO	054410-96-7	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW120221\
Data File : VW021052.D
Acq On : 02 Dec 2021 17:13
Operator : SY/VA
Sample : M4886-18
Misc : 5.00g/10.0mL/MSVOA_W/SOIL
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
EX897

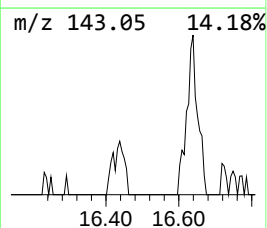
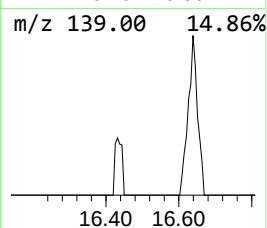
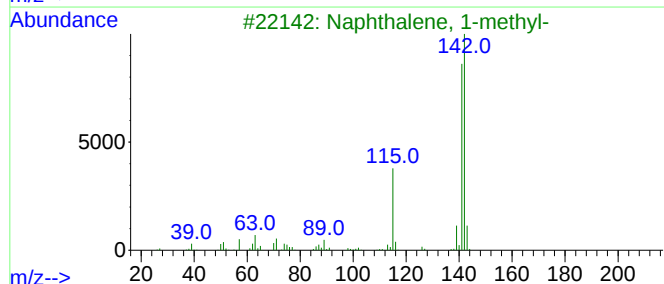
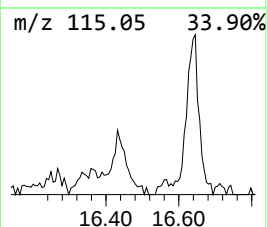
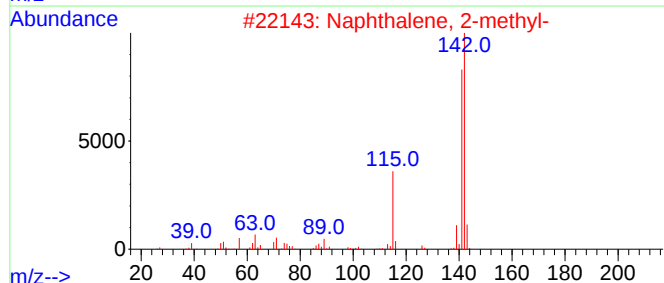
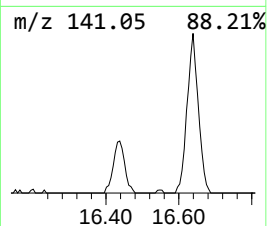
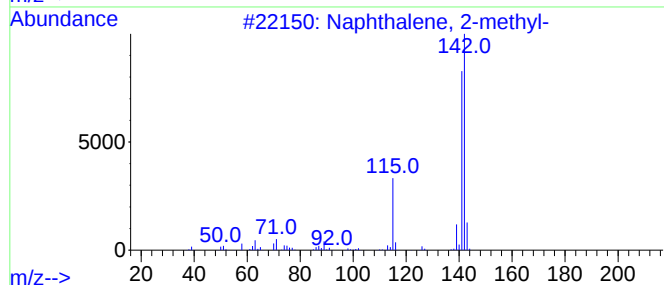
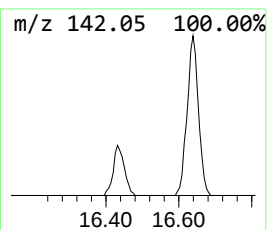
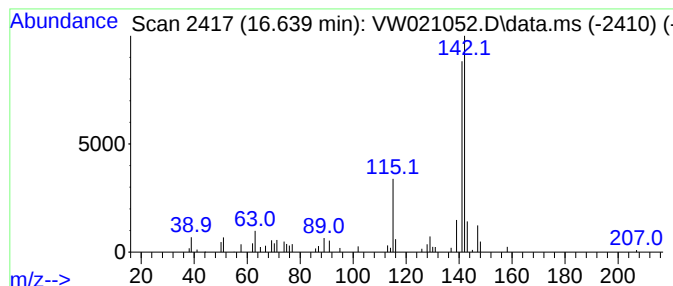
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Quant Title : SFAM01.0

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

Peak Number 9 Naphthalene, 2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.639	2.35 ug/L	30488	1,4-Dichlorobenzene-d4	13.554

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2-methyl-	142	C11H10	000091-57-6	94
2			Naphthalene, 2-methyl-	142	C11H10	000091-57-6	93
3			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	93
4			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	93
5			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	93



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW120221\
Data File : VW021052.D
Acq On : 02 Dec 2021 17:13
Operator : SY/VA
Sample : M4886-18
Misc : 5.00g/10.0mL/MSVOA_W/SOIL
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
EX897

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM111521SMA.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
Naphthalene, 2,...	12.878	3.1	ug/L	39984	3	13.554	324033	25.0
Naphthalene, 1,...	13.292	3.5	ug/L	44764	3	13.554	324033	25.0
Acenaphthene	14.950	11.1	ug/L	143181	3	13.554	324033	25.0
unknown-01	15.066	1.3	ug/L	17083	3	13.554	324033	25.0
Azulene	15.359	2.1	ug/L	27196	3	13.554	324033	25.0
Dibenzofuran	15.633	2.5	ug/L	33103	3	13.554	324033	25.0
Naphthalene, 2-...	16.639	2.4	ug/L	30488	3	13.554	324033	25.0