

Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW122221\
 Data File : VW021569.D
 Acq On : 22 Dec 2021 17:41
 Operator : SY/VA
 Sample : VIBLK562
 Misc : 5.00g/10.0mL/MSVOA_W/SOIL
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 VIBLK562

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\SFAMWLM120921SMA.M
 Title : SFAM01.0

Signal : TIC: VW021569.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.148	37	40	45	rBV2	436	657	0.14%	0.014%
2	2.349	66	73	81	rBV	35537	60915	12.93%	1.304%
3	2.532	101	103	104	rBV	238	132	0.03%	0.003%
4	2.648	120	122	124	rBV	227	145	0.03%	0.003%
5	2.879	153	160	168	rBV	22313	45525	9.66%	0.974%
6	3.379	236	242	243	rBV2	461	635	0.13%	0.014%
7	3.586	273	276	278	rBV	124	171	0.04%	0.004%
8	3.617	278	281	284	rVV	112	186	0.04%	0.004%
9	3.782	306	308	311	rBV	92	137	0.03%	0.003%
10	4.013	335	346	362	rVV	52955	152079	32.28%	3.255%
11	4.129	362	365	366	rVV	204	217	0.05%	0.005%
12	4.147	366	368	371	rVV2	317	368	0.08%	0.008%
13	4.458	409	419	422	rBV3	552	1557	0.33%	0.033%
14	4.659	447	452	453	rBV	86	145	0.03%	0.003%
15	4.745	464	466	469	rVV	119	163	0.03%	0.003%
16	4.909	483	493	504	rVV2	5512	17593	3.73%	0.377%
17	5.098	521	524	525	rBV2	170	216	0.05%	0.005%
18	5.117	525	527	532	rVV	305	415	0.09%	0.009%
19	5.184	536	538	543	rVV2	103	155	0.03%	0.003%
20	5.568	598	601	605	rBB	127	156	0.03%	0.003%
21	5.647	613	614	617	rBV	118	124	0.03%	0.003%
22	5.751	628	631	636	rVB	243	283	0.06%	0.006%
23	5.806	636	640	642	rBB	115	147	0.03%	0.003%
24	5.842	643	646	648	rBV	130	162	0.03%	0.003%
25	5.940	655	662	671	rBB5	1144	2963	0.63%	0.063%
26	6.074	682	684	689	rVB	127	172	0.04%	0.004%
27	6.281	716	718	720	rBV	123	140	0.03%	0.003%
28	6.598	767	770	772	rBB	154	135	0.03%	0.003%
29	6.860	809	813	814	rBV	154	178	0.04%	0.004%
30	6.872	814	815	818	rBV	115	123	0.03%	0.003%
31	7.080	840	849	856	rBV	12184	33422	7.09%	0.715%
32	7.330	888	890	893	rVV	124	193	0.04%	0.004%
33	7.647	930	942	954	rBB	79334	198110	42.04%	4.240%
34	7.763	958	961	963	rBB	135	139	0.03%	0.003%
35	7.927	985	988	992	rBB	81	139	0.03%	0.003%
36	8.092	1009	1015	1016	rBB	129	206	0.04%	0.004%

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Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM120921SMA.M
 Title : SFAM01.0

37	8.275	1034	1045	1059	rBV2	157631	471194	100.00%	10.085%
38	8.403	1063	1066	1068	rBV2	328	412	0.09%	0.009%
39	8.488	1078	1080	1084	rVB2	464	540	0.11%	0.012%
40	8.543	1088	1089	1092	rBV	155	173	0.04%	0.004%
41	8.598	1097	1098	1102	rVB2	180	151	0.03%	0.003%
42	8.628	1102	1103	1106	rVB2	152	124	0.03%	0.003%
43	8.842	1128	1138	1147	rBV	142194	276291	58.64%	5.914%
44	8.909	1147	1149	1150	rVV	202	149	0.03%	0.003%
45	9.079	1173	1177	1180	rBV2	533	724	0.15%	0.015%
46	9.274	1200	1209	1218	rVV	105253	213587	45.33%	4.572%
47	9.354	1218	1222	1224	rVV2	292	437	0.09%	0.009%
48	9.384	1224	1227	1229	rVB	115	155	0.03%	0.003%
49	9.506	1245	1247	1249	rVB	171	132	0.03%	0.003%
50	9.622	1263	1266	1268	rVB	127	148	0.03%	0.003%
51	9.659	1268	1272	1273	rBV	328	350	0.07%	0.007%
52	9.890	1306	1310	1314	rBB	333	459	0.10%	0.010%
53	9.982	1323	1325	1326	rBV	271	185	0.04%	0.004%
54	10.037	1326	1334	1342	rVV	58507	105254	22.34%	2.253%
55	10.091	1342	1343	1346	rVB	243	151	0.03%	0.003%
56	10.213	1358	1363	1365	rBV2	506	625	0.13%	0.013%
57	10.323	1373	1381	1388	rBV	236994	422056	89.57%	9.033%
58	10.384	1388	1391	1400	rVB	3347	5893	1.25%	0.126%
59	10.579	1416	1423	1430	rBV	37459	63430	13.46%	1.358%
60	10.652	1433	1435	1436	rVB	205	135	0.03%	0.003%
61	10.701	1436	1443	1449	rBV	17300	30552	6.48%	0.654%
62	10.780	1455	1456	1459	rVB3	358	311	0.07%	0.007%
63	10.811	1459	1461	1462	rBV	294	263	0.06%	0.006%
64	10.841	1464	1466	1467	rBV2	521	376	0.08%	0.008%
65	10.878	1470	1472	1473	rBV	328	208	0.04%	0.004%
66	10.921	1473	1479	1487	rBV	50886	89394	18.97%	1.913%
67	11.061	1499	1502	1507	rBV2	6126	9636	2.05%	0.206%
68	11.366	1550	1552	1553	rBV2	1082	947	0.20%	0.020%
69	11.439	1559	1564	1570	rVB	10150	18453	3.92%	0.395%
70	11.628	1588	1595	1603	rVB	212925	361086	76.63%	7.729%
71	11.713	1608	1609	1613	rVB2	451	344	0.07%	0.007%
72	11.780	1618	1620	1622	rBV	345	315	0.07%	0.007%
73	11.835	1626	1629	1633	rBV4	670	1119	0.24%	0.024%
74	11.908	1638	1641	1642	rBV3	561	603	0.13%	0.013%
75	12.024	1656	1660	1663	rVB4	852	898	0.19%	0.019%
76	12.128	1675	1677	1680	rBV2	433	469	0.10%	0.010%

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 Title : SFAM01.0

77	12.280	1690	1702	1703	rBV9	2688	7397	1.57%	0.158%
78	12.603	1749	1755	1761	rVB2	59279	97354	20.66%	2.084%
79	12.689	1763	1769	1776	rVB2	90215	148742	31.57%	3.184%
80	12.865	1776	1798	1821	rBV2	45346	297766	63.19%	6.373%
81	13.310	1858	1871	1884	rBV6	74334	356311	75.62%	7.626%
82	13.554	1905	1911	1917	rBV	268416	417458	88.60%	8.935%
83	13.853	1952	1960	1970	rBV	256934	468514	99.43%	10.028%
84	14.066	1990	1995	2010	rVB2	23864	49096	10.42%	1.051%
85	14.298	2024	2033	2045	rVB2	14603	51056	10.84%	1.093%
86	14.725	2099	2103	2104	rBV3	3089	4027	0.85%	0.086%
87	15.066	2153	2159	2168	rBV3	7587	21972	4.66%	0.470%
88	15.359	2199	2207	2212	rBV	16193	31987	6.79%	0.685%
89	15.432	2214	2219	2223	rBV	11913	20807	4.42%	0.445%
90	15.621	2241	2250	2263	rBV6	7646	27999	5.94%	0.599%
91	15.944	2294	2303	2311	rVB8	4205	13128	2.79%	0.281%
92	16.072	2320	2324	2325	rBV3	777	1034	0.22%	0.022%
93	16.212	2345	2347	2348	rVB3	674	365	0.08%	0.008%
94	16.231	2348	2350	2351	rVB	219	144	0.03%	0.003%
95	16.243	2351	2352	2354	rBV	418	355	0.08%	0.008%
96	16.352	2362	2370	2371	rBV6	2535	4660	0.99%	0.100%
97	16.438	2377	2384	2395	rVB4	6256	17308	3.67%	0.370%
98	16.554	2399	2403	2405	rVB3	444	561	0.12%	0.012%
99	16.639	2407	2417	2428	rBV2	15731	37889	8.04%	0.811%
100	16.730	2430	2432	2434	rBV2	282	271	0.06%	0.006%

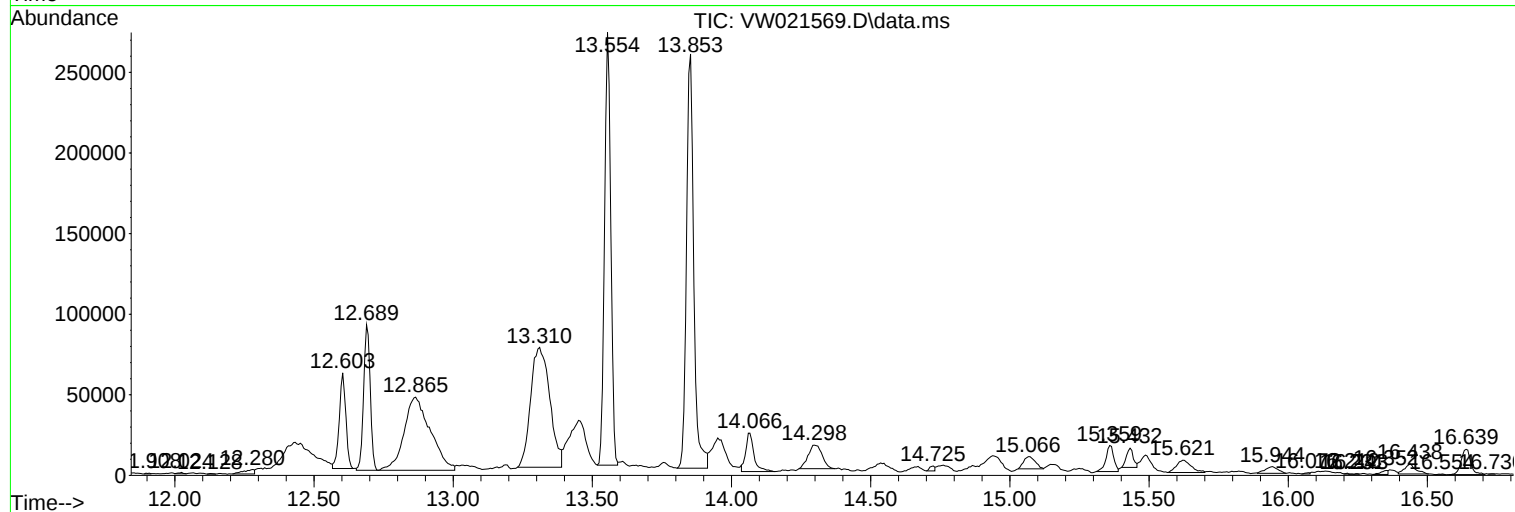
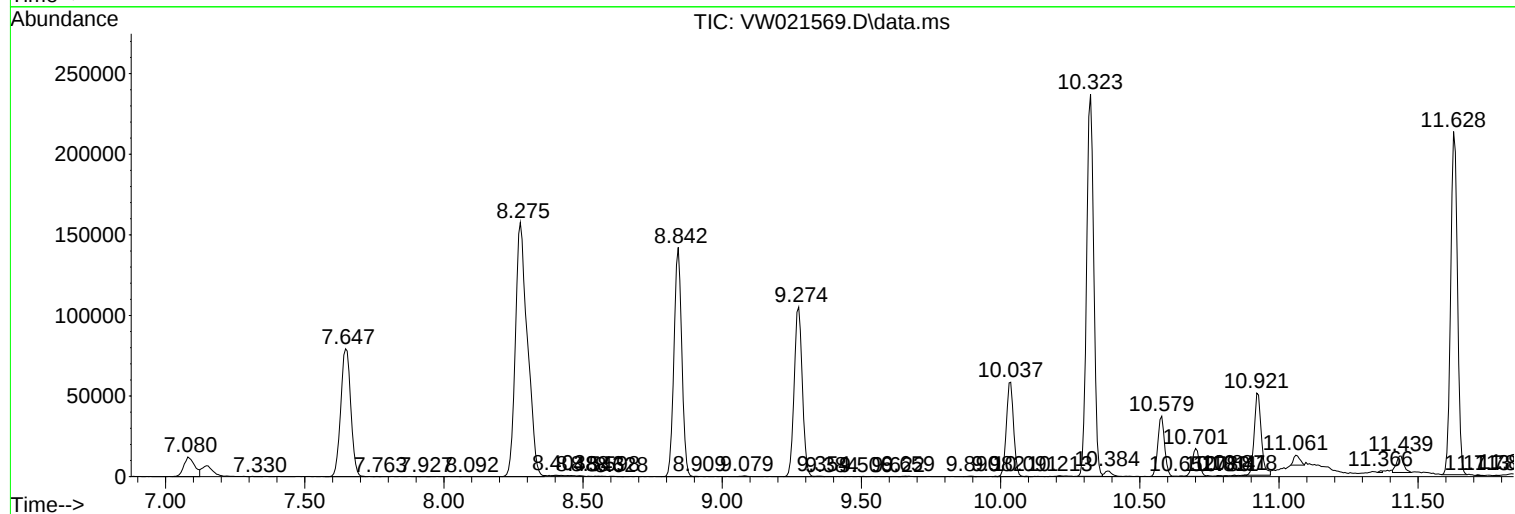
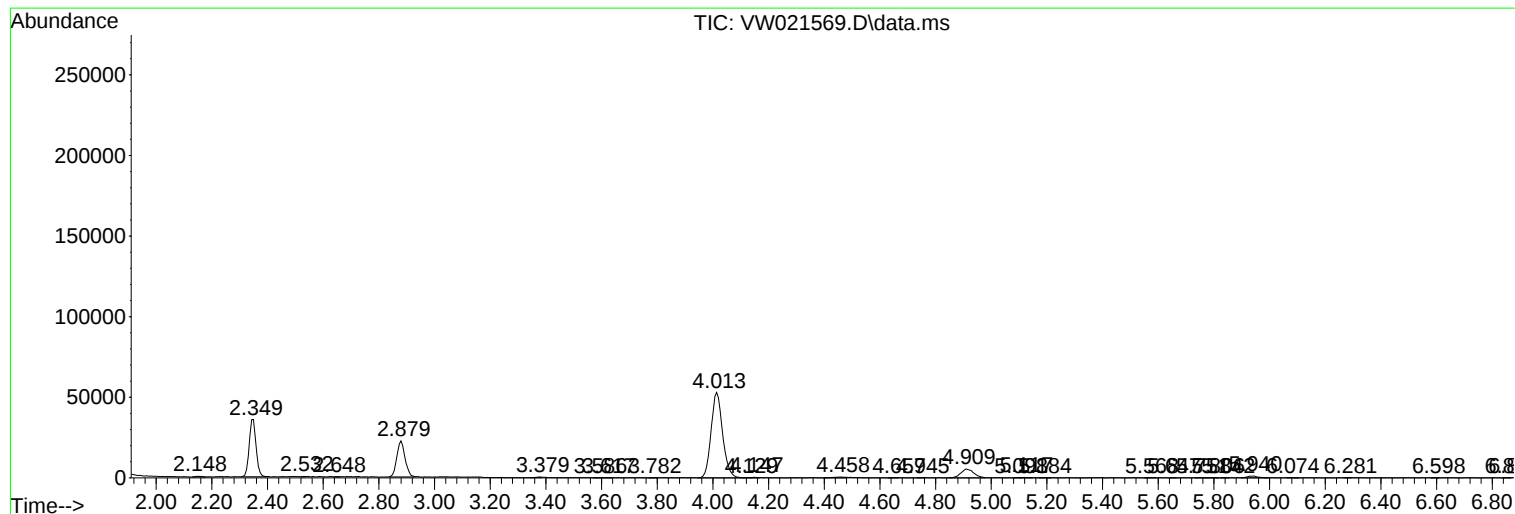
Sum of corrected areas: 4672133

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Instrument :
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM120921SMA.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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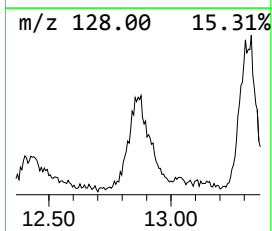
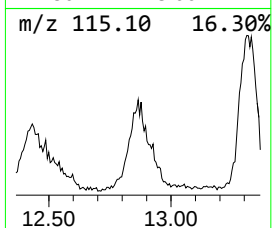
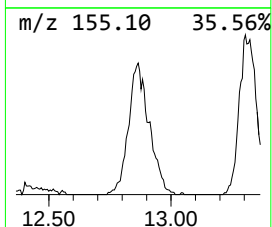
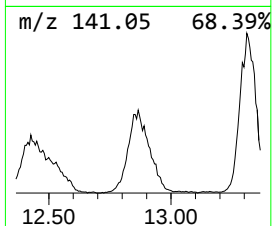
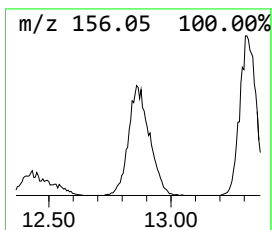
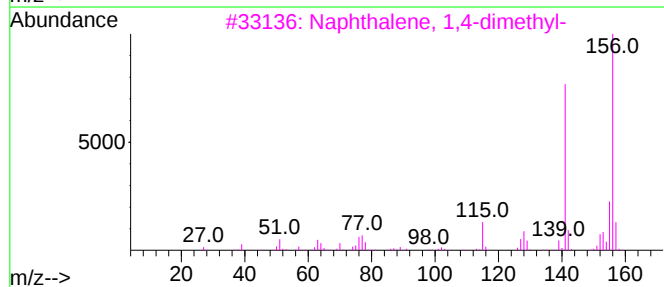
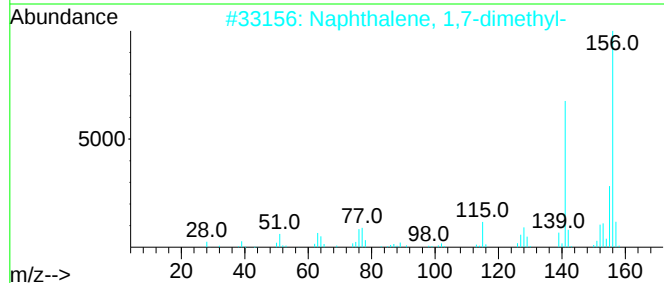
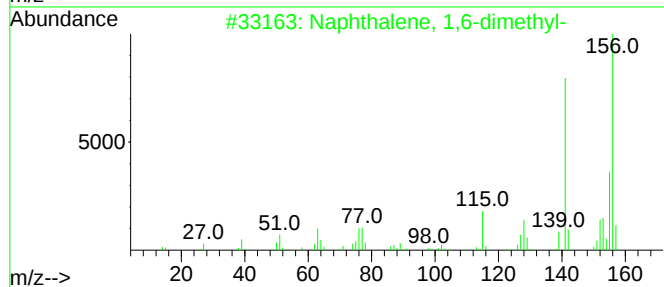
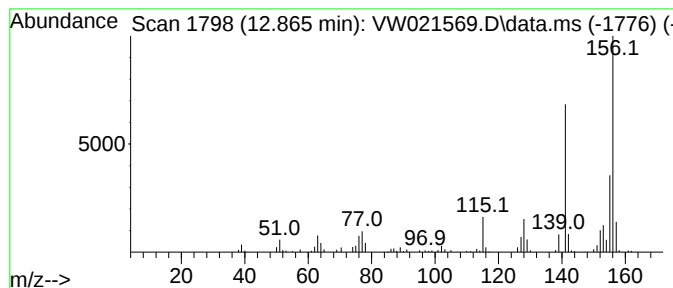
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM120921SMA.M
Quant Title : SFAM01.0

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

Peak Number 6 Naphthalene, 1,6-dimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.865	17.83 ug/L	297766	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	97
2			Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	97
3			Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	97
4			Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	97
5			Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	96



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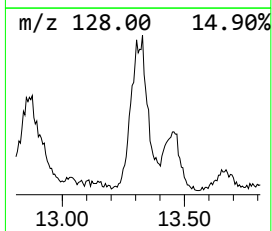
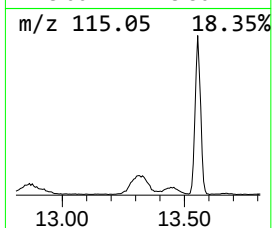
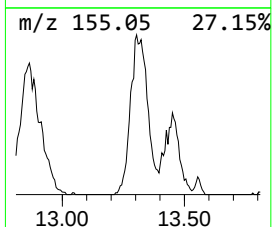
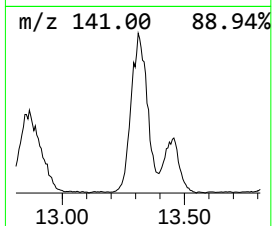
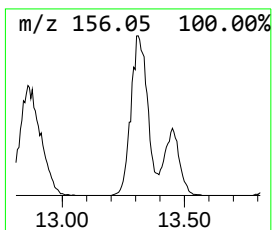
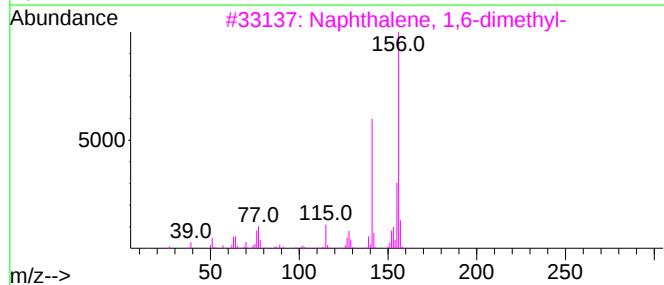
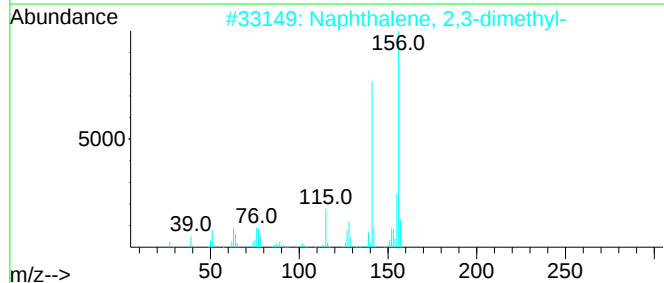
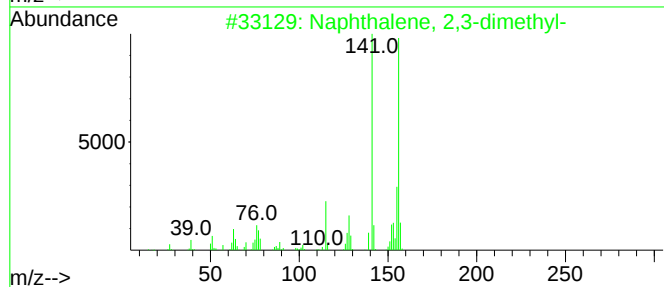
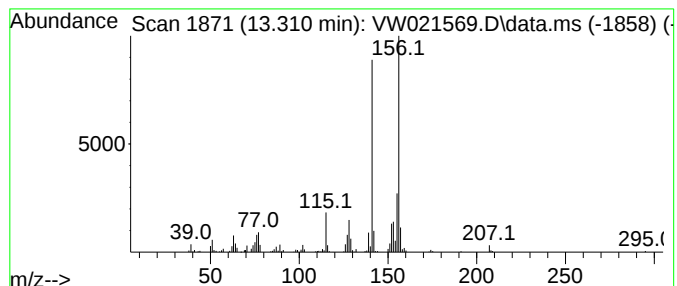
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM120921SMA.M
Quant Title : SFAM01.0

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

Peak Number 7 Naphthalene, 2,3-dimethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.310	21.34 ug/L	356311	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	98
2			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
3			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	97
4			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	97
5			Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	97



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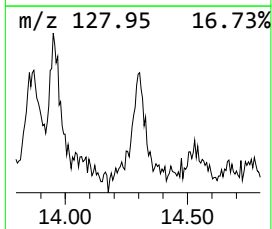
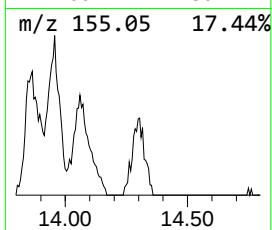
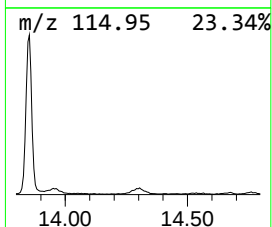
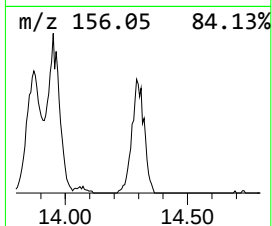
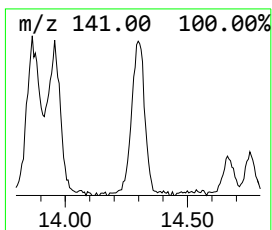
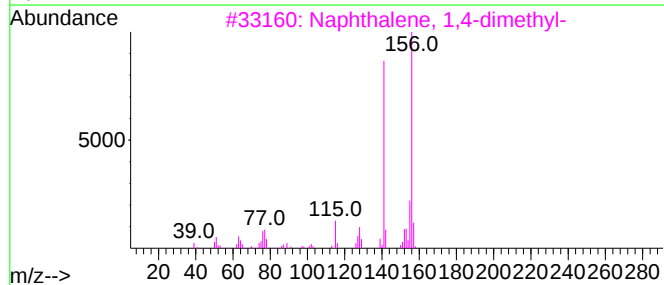
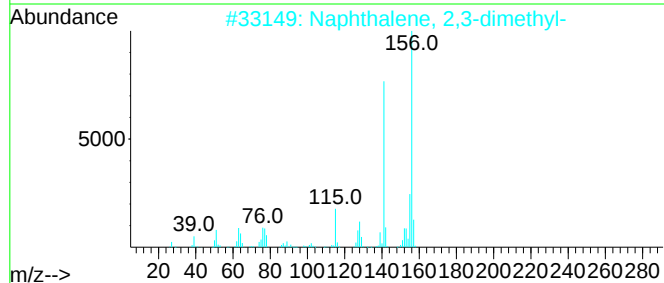
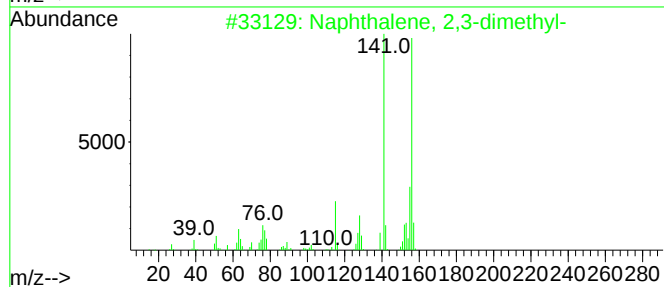
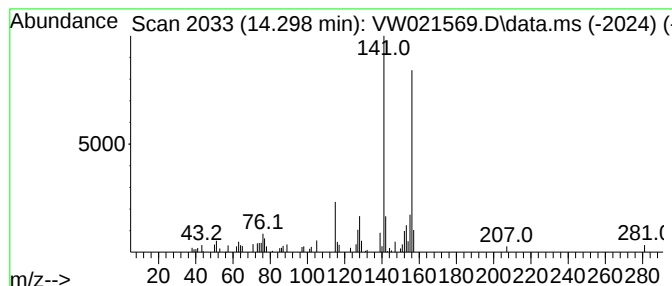
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM120921SMA.M
Quant Title : SFAM01.0

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

Peak Number 9 Naphthalene, 1,4-dimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.298	3.06 ug/L	51056	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	96
2			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	96
3			Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	96
4			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	95
5			Naphthalene, 1,8-dimethyl-	156	C12H12	000569-41-5	95



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW122221\
Data File : VW021569.D
Acq On : 22 Dec 2021 17:41
Operator : SY/VA
Sample : VIBLK562
Misc : 5.00g/10.0mL/MSVOA_W/SOIL
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
VIBLK562

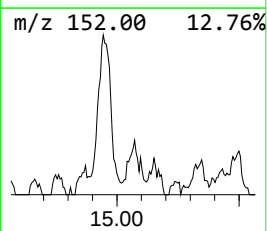
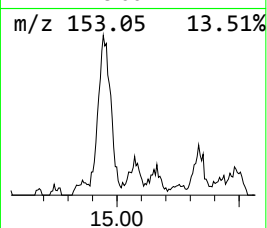
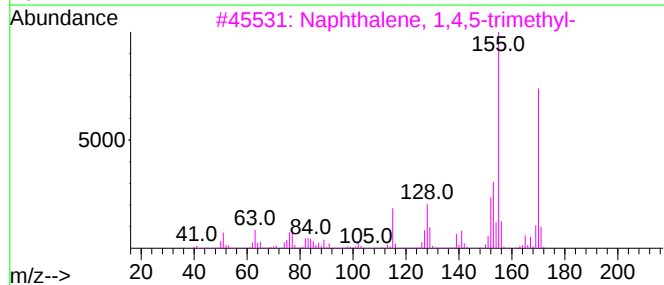
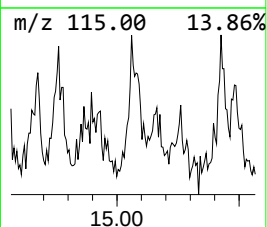
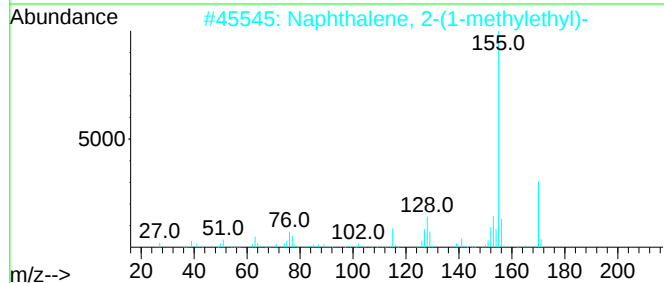
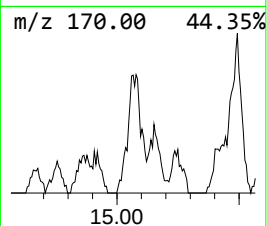
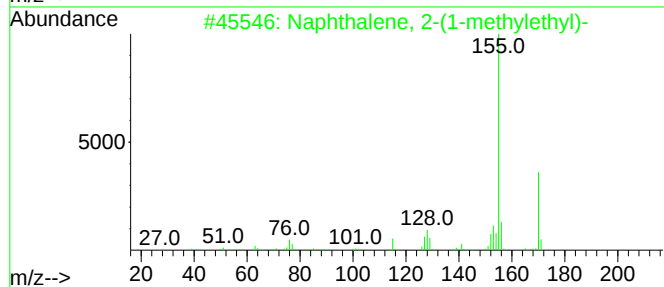
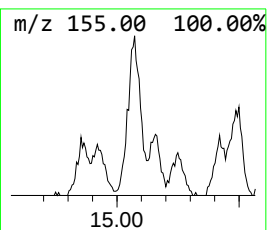
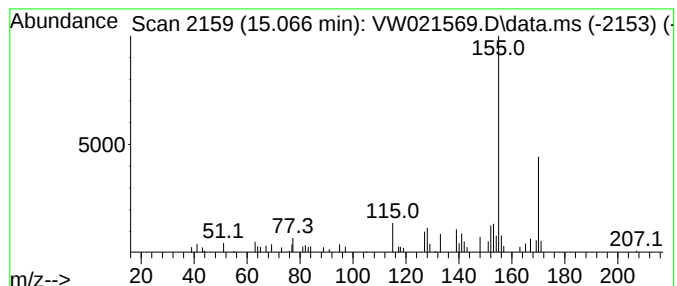
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Quant Title : SFAM01.0

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

Peak Number 10 Naphthalene, 2-(1-methyleth... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.066	1.32 ug/L	21972	1,4-Dichlorobenzene-d4	13.554

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2-(1-methylethyl)-	170	C13H14	002027-17-0	87
2			Naphthalene, 2-(1-methylethyl)-	170	C13H14	002027-17-0	87
3			Naphthalene, 1,4,5-trimethyl-	170	C13H14	002131-41-1	64
4			3-(2-Methyl-propenyl)-1H-indene	170	C13H14	1000187-78-5	64
5			2-Naphthyl methyl ketone	170	C12H10O	000093-08-3	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW122221\
Data File : VW021569.D
Acq On : 22 Dec 2021 17:41
Operator : SY/VA
Sample : VIBLK562
Misc : 5.00g/10.0mL/MSVOA_W/SOIL
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
VIBLK562

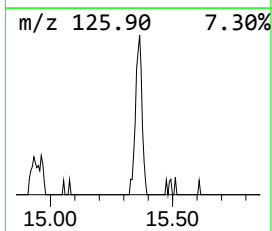
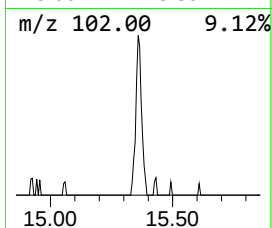
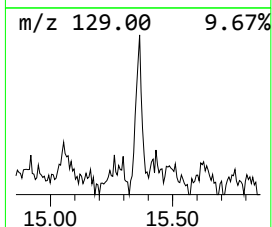
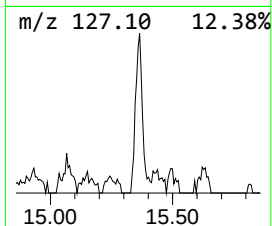
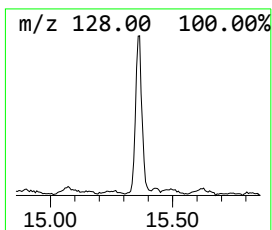
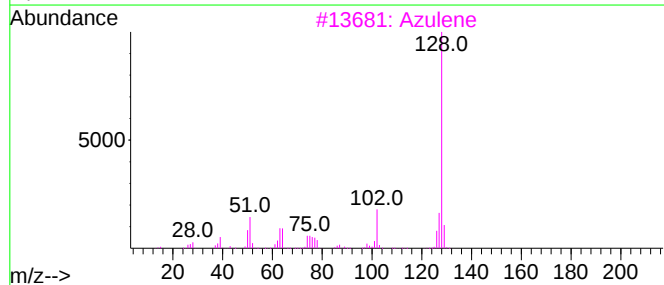
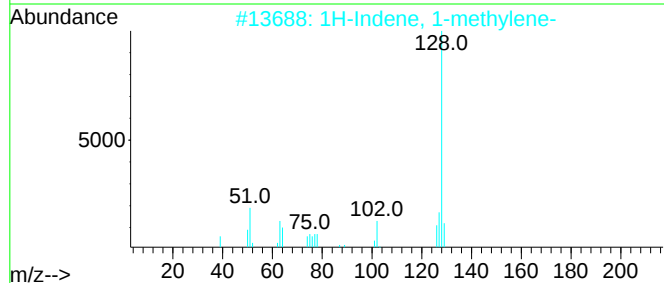
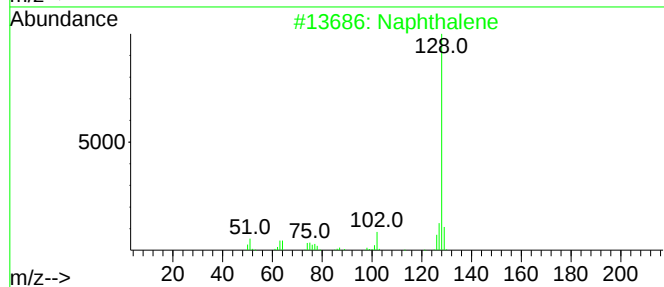
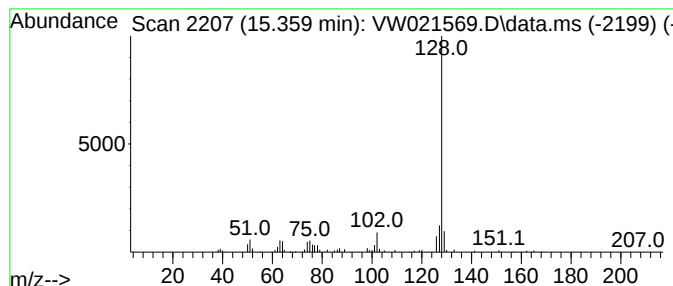
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Quant Title : SFAM01.0

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

Peak Number 11 Azulene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.359	1.92 ug/L	31987	1,4-Dichlorobenzene-d4	13.554

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene	128	C10H8	000091-20-3	95
2			1H-Indene, 1-methylene-	128	C10H8	002471-84-3	94
3			Azulene	128	C10H8	000275-51-4	91
4			Azulene	128	C10H8	000275-51-4	91
5			Naphthalene	128	C10H8	000091-20-3	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW122221\
Data File : VW021569.D
Acq On : 22 Dec 2021 17:41
Operator : SY/VA
Sample : VIBLK562
Misc : 5.00g/10.0mL/MSVOA_W/SOIL
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
VIBLK562

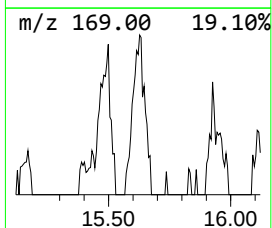
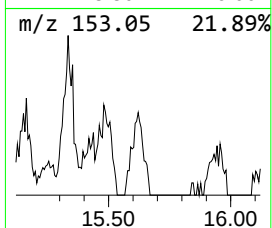
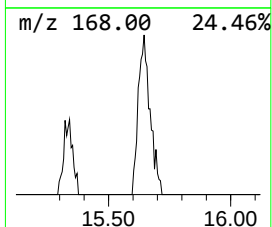
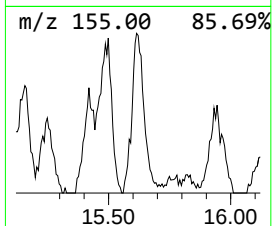
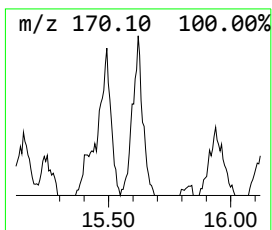
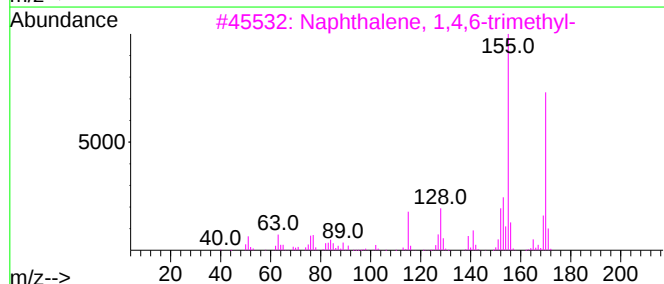
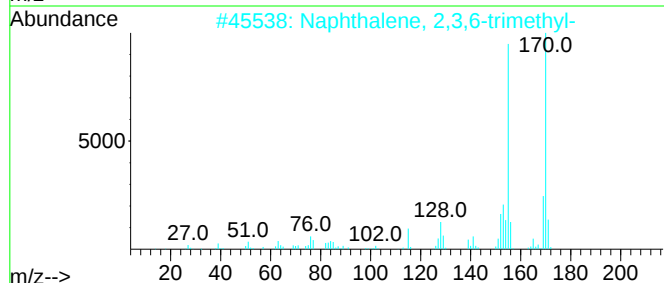
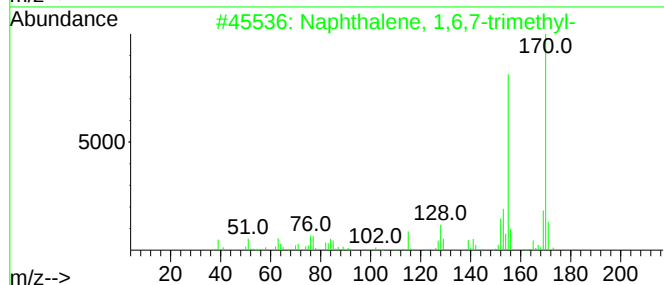
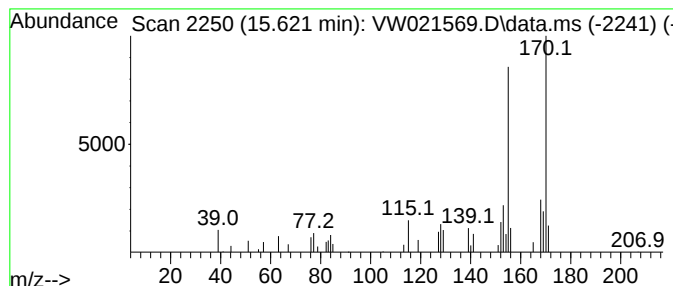
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Quant Title : SFAM01.0

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

Peak Number 12 Naphthalene, 1,6,7-trimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.621	1.68 ug/L	27999	1,4-Dichlorobenzene-d4	13.554

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	96
2			Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	95
3			Naphthalene, 1,4,6-trimethyl-	170	C13H14	002131-42-2	95
4			Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	95
5			Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	95



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW122221\
Data File : VW021569.D
Acq On : 22 Dec 2021 17:41
Operator : SY/VA
Sample : VIBLK562
Misc : 5.00g/10.0mL/MSVOA_W/SOIL
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
VIBLK562

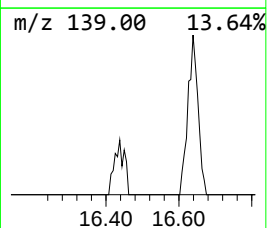
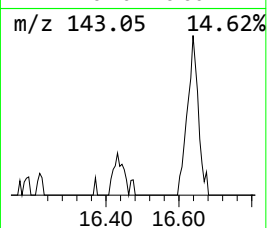
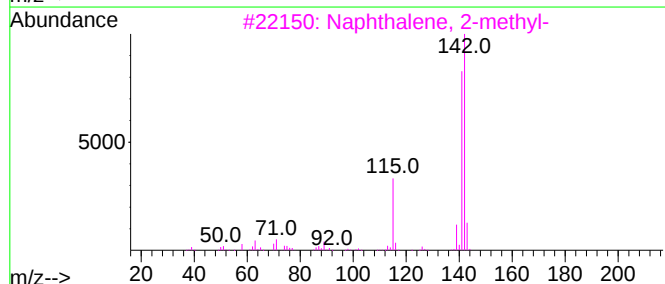
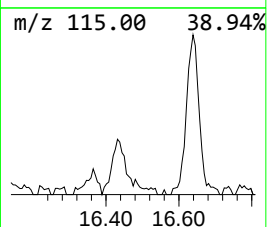
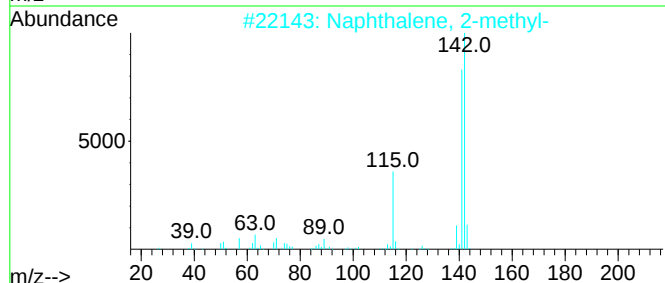
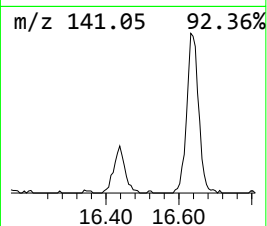
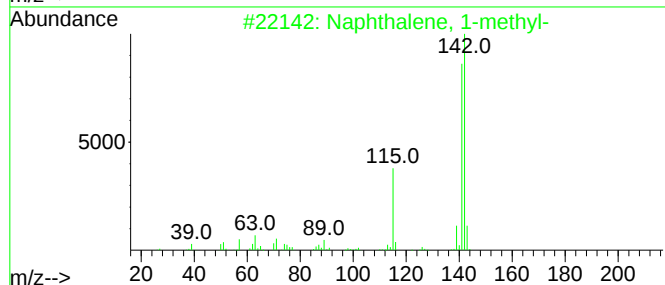
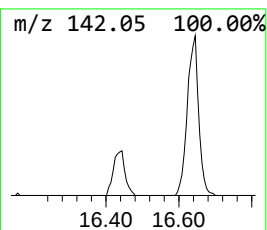
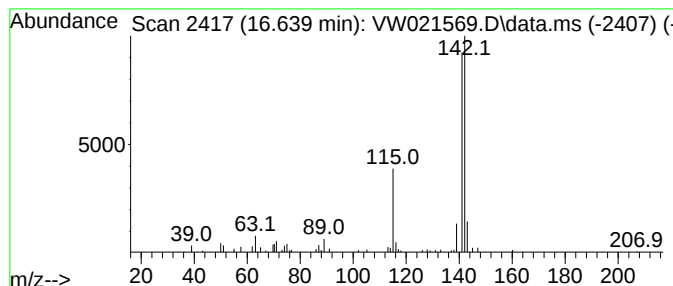
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Quant Title : SFAM01.0

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

Peak Number 13 1,4-Methanonaphthalene, 1,4... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.639	2.27 ug/L	37889	1,4-Dichlorobenzene-d4	13.554

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	97
2			Naphthalene, 2-methyl-	142	C11H10	000091-57-6	97
3			Naphthalene, 2-methyl-	142	C11H10	000091-57-6	95
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\VW122221\
Data File : VW021569.D
Acq On : 22 Dec 2021 17:41
Operator : SY/VA
Sample : VIBLK562
Misc : 5.00g/10.0mL/MSVOA_W/SOIL
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
VIBLK562

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM120921SMA.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, 1,...	12.865	17.8	ug/L	297766	3	13.554	417458	25.0
Naphthalene, 2,...	13.310	21.3	ug/L	356311	3	13.554	417458	25.0
Naphthalene, 1,...	14.298	3.1	ug/L	51056	3	13.554	417458	25.0
Naphthalene, 2-...	15.066	1.3	ug/L	21972	3	13.554	417458	25.0
Azulene	15.359	1.9	ug/L	31987	3	13.554	417458	25.0
Naphthalene, 1,...	15.621	1.7	ug/L	27999	3	13.554	417458	25.0
1,4-Methanonaph...	16.639	2.3	ug/L	37889	3	13.554	417458	25.0