

Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW101121\
 Data File : VW020444.D
 Acq On : 11 Oct 2021 15:52
 Operator : SY/VA
 Sample : VIBLK472
 Misc : 5.00g/10mL/MSVOA_W/SOIL
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 VIBLK472

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\SFAMWLM100821SMA.M
 Title : SFAM01.0

Signal : TIC: VW020444.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.349	67	73	82	rVB	85334	134365	12.10%	1.581%
2	2.885	154	161	172	rVB	39942	80656	7.26%	0.949%
3	3.672	288	290	296	rVB3	275	418	0.04%	0.005%
4	3.751	301	303	305	rVB	127	96	0.01%	0.001%
5	3.782	305	308	311	rBV	81	110	0.01%	0.001%
6	4.013	335	346	363	rBV	124466	360216	32.44%	4.238%
7	4.129	363	365	367	rVV	264	208	0.02%	0.002%
8	4.269	387	388	390	rVB2	124	88	0.01%	0.001%
9	4.690	455	457	459	rBV	148	108	0.01%	0.001%
10	4.745	461	466	467	rBV2	87	129	0.01%	0.002%
11	4.915	481	494	509	rBV2	34092	111875	10.07%	1.316%
12	5.324	560	561	564	rBV	121	102	0.01%	0.001%
13	5.409	572	575	579	rBV2	104	144	0.01%	0.002%
14	5.604	605	607	611	rVB2	86	130	0.01%	0.002%
15	5.671	616	618	619	rVB2	124	84	0.01%	0.001%
16	5.690	619	621	622	rBV	118	84	0.01%	0.001%
17	5.946	651	663	673	rVV2	8433	24949	2.25%	0.294%
18	6.293	719	720	723	rBV	136	140	0.01%	0.002%
19	6.366	728	732	735	rVB2	133	224	0.02%	0.003%
20	6.397	735	737	739	rBV	93	97	0.01%	0.001%
21	6.604	769	771	773	rVB	197	104	0.01%	0.001%
22	6.830	805	808	809	rBV2	172	117	0.01%	0.001%
23	7.086	840	850	871	rBV	31399	89385	8.05%	1.052%
24	7.244	875	876	878	rVV2	139	118	0.01%	0.001%
25	7.360	892	895	898	rVB2	142	151	0.01%	0.002%
26	7.415	901	904	907	rVB2	90	127	0.01%	0.001%
27	7.446	907	909	911	rVB2	117	101	0.01%	0.001%
28	7.543	921	925	931	rVB3	504	1005	0.09%	0.012%
29	7.647	931	942	957	rBV	185181	452478	40.75%	5.323%
30	7.872	976	979	980	rBV	157	144	0.01%	0.002%
31	7.945	987	991	995	rBV4	286	477	0.04%	0.006%
32	8.012	1001	1002	1004	rBV2	92	87	0.01%	0.001%
33	8.043	1004	1007	1009	rVB	103	107	0.01%	0.001%
34	8.067	1009	1011	1013	rBV2	80	94	0.01%	0.001%
35	8.122	1018	1020	1024	rVB2	118	108	0.01%	0.001%
36	8.177	1024	1029	1032	rVB	146	168	0.02%	0.002%

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Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM100821SMA.M
 Title : SFAM01.0

37	8.275	1035	1045	1060	rBV2	374072	1110486	100.00%	13.064%
38	8.689	1108	1113	1117	rBV3	267	426	0.04%	0.005%
39	8.842	1129	1138	1153	rBV	379233	742639	66.88%	8.737%
40	8.957	1156	1157	1159	rVB2	165	86	0.01%	0.001%
41	9.079	1172	1177	1182	rBV3	684	1090	0.10%	0.013%
42	9.201	1193	1197	1198	rBV2	110	97	0.01%	0.001%
43	9.274	1200	1209	1219	rBV	269118	531137	47.83%	6.249%
44	9.494	1239	1245	1249	rBV5	251	499	0.04%	0.006%
45	9.561	1255	1256	1258	rVB2	143	101	0.01%	0.001%
46	9.591	1258	1261	1264	rBV2	147	185	0.02%	0.002%
47	9.628	1264	1267	1269	rVV	84	104	0.01%	0.001%
48	9.665	1269	1273	1276	rVV3	349	476	0.04%	0.006%
49	9.695	1276	1278	1279	rVV2	158	114	0.01%	0.001%
50	9.713	1279	1281	1283	rVB	179	165	0.01%	0.002%
51	9.744	1283	1286	1287	rBV2	147	138	0.01%	0.002%
52	9.756	1287	1288	1291	rBV2	114	108	0.01%	0.001%
53	9.817	1296	1298	1300	rBV2	140	143	0.01%	0.002%
54	9.866	1302	1306	1308	rBV2	96	143	0.01%	0.002%
55	9.896	1308	1311	1313	rVB3	91	119	0.01%	0.001%
56	9.939	1317	1318	1321	rVB2	156	134	0.01%	0.002%
57	9.976	1321	1324	1326	rBV3	154	193	0.02%	0.002%
58	10.037	1326	1334	1348	rBV	148037	265047	23.87%	3.118%
59	10.183	1356	1358	1359	rBV2	214	180	0.02%	0.002%
60	10.225	1363	1365	1369	rVV2	200	274	0.02%	0.003%
61	10.323	1373	1381	1388	rVV	553890	953890	85.90%	11.222%
62	10.390	1388	1392	1403	rVB	8232	14964	1.35%	0.176%
63	10.469	1403	1405	1408	rVB2	174	137	0.01%	0.002%
64	10.500	1408	1410	1414	rVB3	359	353	0.03%	0.004%
65	10.530	1414	1415	1416	rBV	147	102	0.01%	0.001%
66	10.579	1416	1423	1432	rBV	95136	162815	14.66%	1.915%
67	10.640	1432	1433	1435	rBV	314	173	0.02%	0.002%
68	10.664	1435	1437	1439	rBV3	271	192	0.02%	0.002%
69	10.707	1439	1444	1449	rVB	3041	5495	0.49%	0.065%
70	10.780	1454	1456	1458	rBV2	157	134	0.01%	0.002%
71	10.817	1460	1462	1463	rBV2	219	148	0.01%	0.002%
72	10.927	1473	1480	1494	rBV	122503	208197	18.75%	2.449%
73	11.024	1494	1496	1498	rVB3	345	281	0.03%	0.003%
74	11.164	1517	1519	1520	rBV2	275	188	0.02%	0.002%
75	11.231	1524	1530	1533	rBV3	754	1290	0.12%	0.015%
76	11.347	1543	1549	1552	rBV5	1202	2187	0.20%	0.026%

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 Title : SFAM01.0

77	11.555	1581	1583	1586	rVB4	380	281	0.03%	0.003%
78	11.634	1588	1596	1605	rVV	517395	899837	81.03%	10.586%
79	11.841	1624	1630	1633	rBV6	1137	2325	0.21%	0.027%
80	12.000	1654	1656	1659	rVB4	457	445	0.04%	0.005%
81	12.036	1659	1662	1665	rBV3	414	633	0.06%	0.007%
82	12.079	1668	1669	1672	rBV3	255	197	0.02%	0.002%
83	12.121	1673	1676	1677	rBV2	623	612	0.06%	0.007%
84	12.164	1680	1683	1685	rBV3	779	1000	0.09%	0.012%
85	12.262	1692	1699	1700	rBV7	2376	4879	0.44%	0.057%
86	12.591	1749	1753	1761	rVB4	9567	19934	1.80%	0.235%
87	12.695	1763	1770	1776	rBV	183608	312026	28.10%	3.671%
88	13.134	1838	1842	1845	rBV5	4526	7010	0.63%	0.082%
89	13.201	1849	1853	1854	rBV2	5612	7918	0.71%	0.093%
90	13.554	1905	1911	1918	rBV	546902	900418	81.08%	10.593%
91	13.853	1953	1960	1967	rBV	480624	792667	71.38%	9.325%
92	14.066	1992	1995	1999	rBV	14528	20995	1.89%	0.247%
93	14.456	2056	2059	2062	rVB	9762	12576	1.13%	0.148%
94	14.493	2062	2065	2069	rBV5	6725	8768	0.79%	0.103%
95	14.694	2094	2098	2100	rBV3	4509	8052	0.73%	0.095%
96	14.798	2110	2115	2119	rBV3	11053	26103	2.35%	0.307%
97	15.322	2196	2201	2206	rBV2	15547	28384	2.56%	0.334%
98	16.279	2354	2358	2363	rBV4	9447	17977	1.62%	0.211%
99	16.438	2379	2384	2396	rVB	37071	90389	8.14%	1.063%
100	16.639	2412	2417	2429	rVB2	27665	73036	6.58%	0.859%

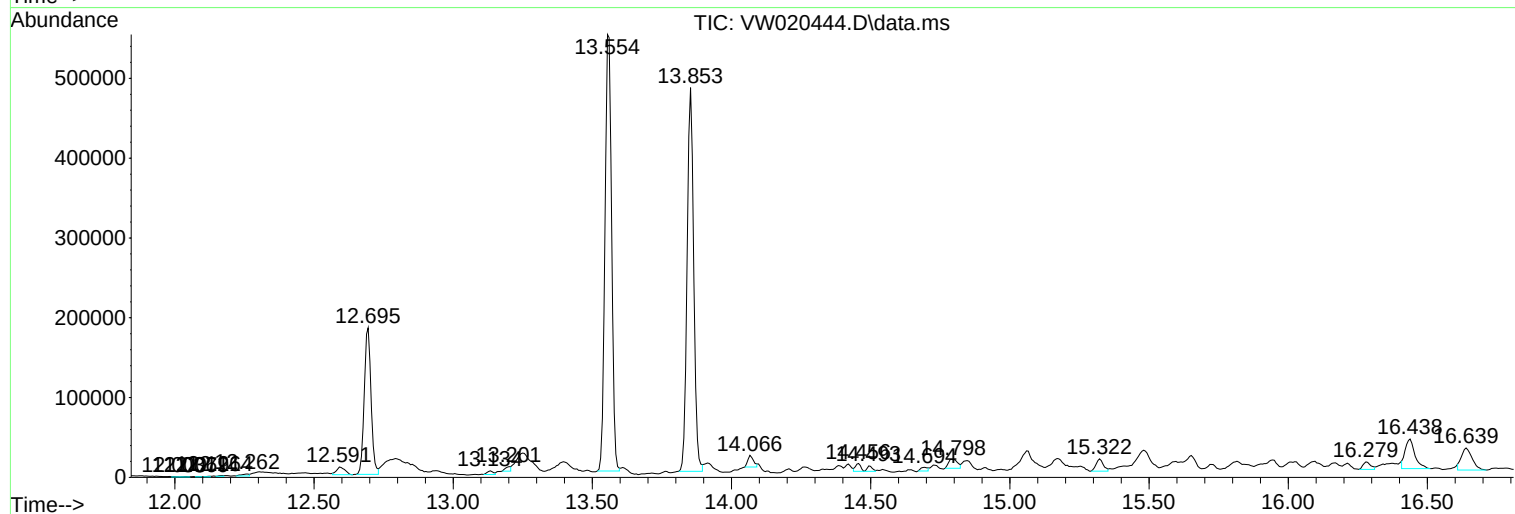
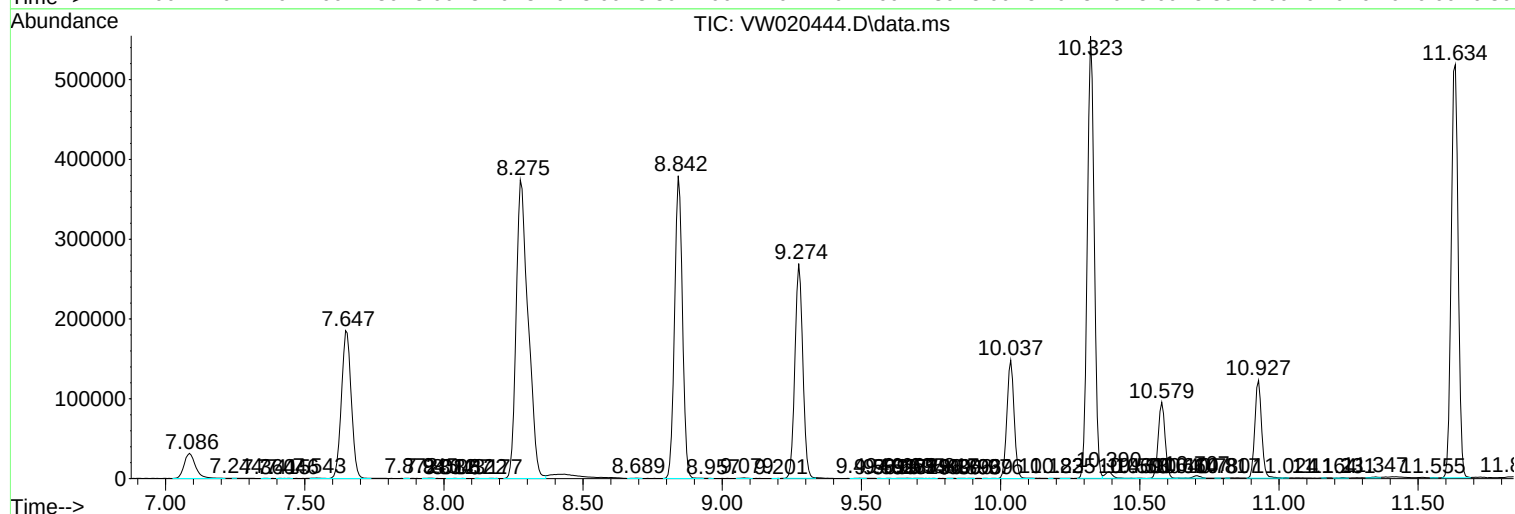
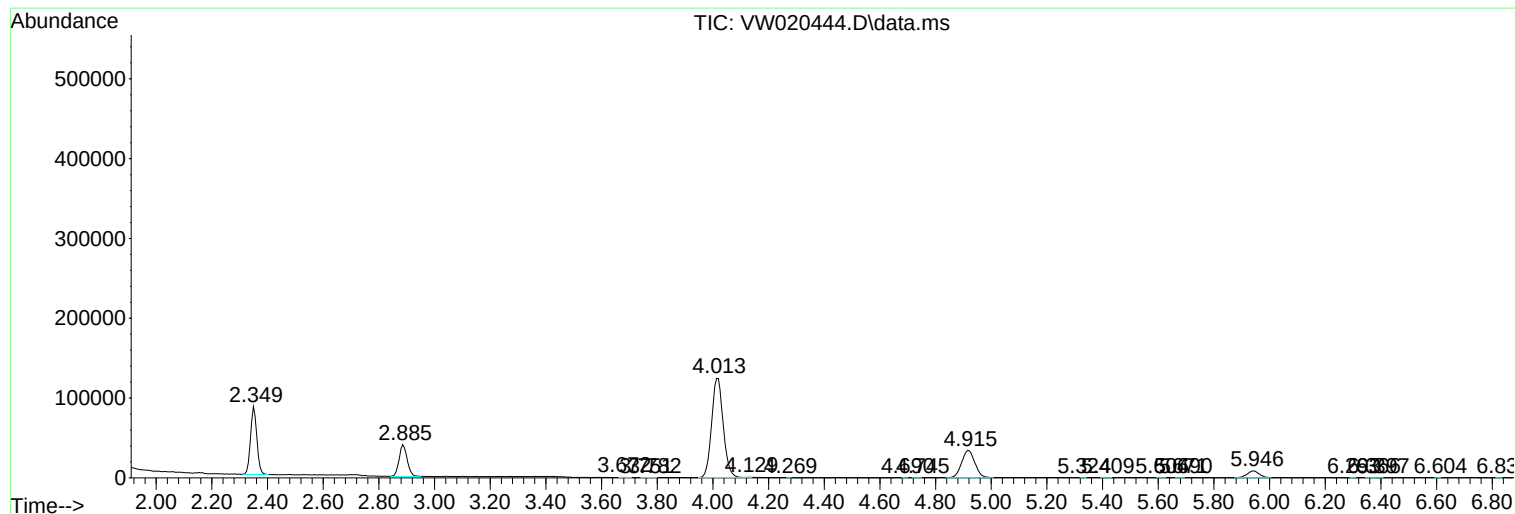
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Instrument :
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM100821SMA.M
Quant Title : SFAM01.0

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



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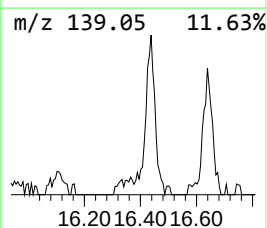
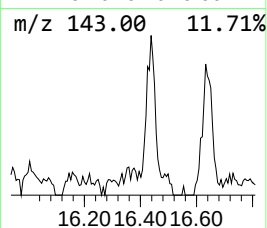
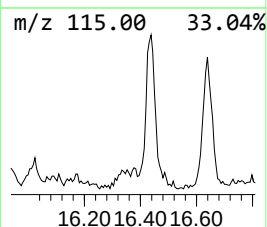
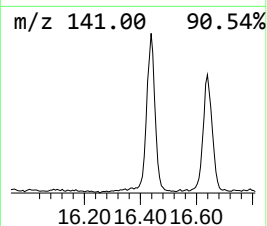
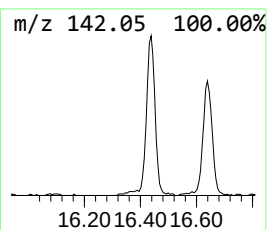
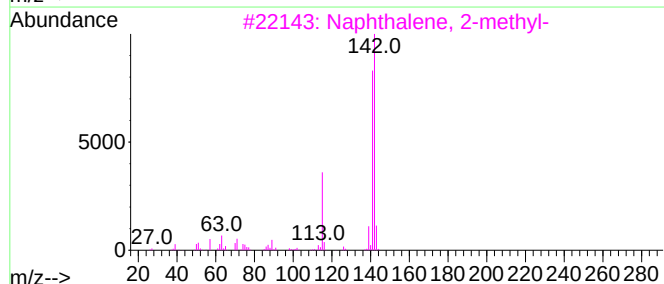
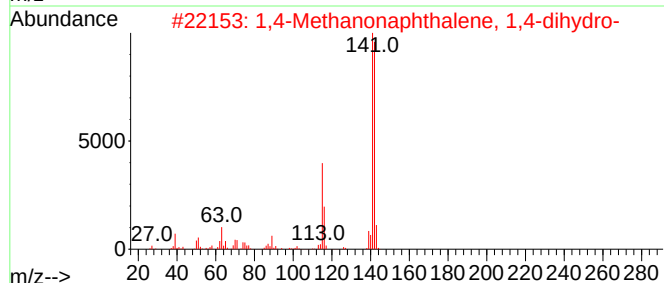
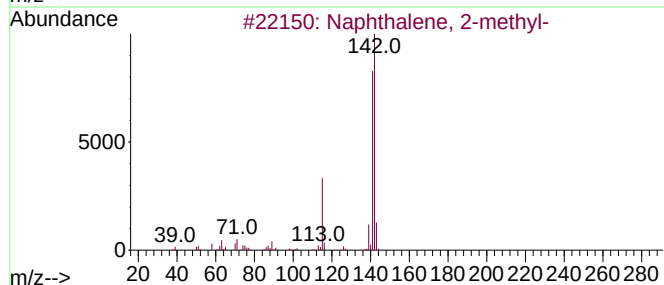
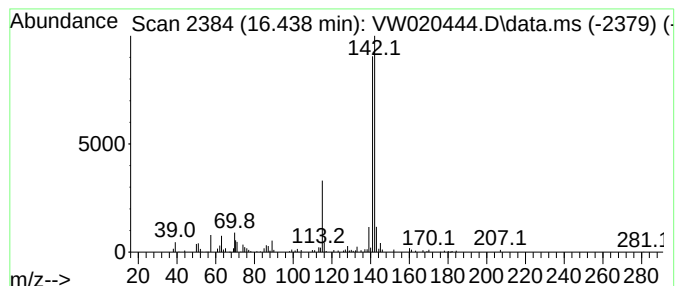
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM100821SMA.M
Quant Title : SFAM01.0

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.438	2.51 ug/L	90389	1,4-Dichlorobenzene-d4	13.560

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



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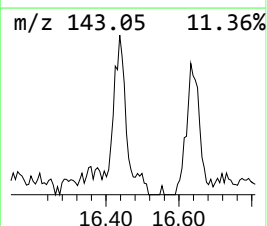
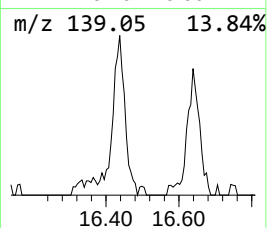
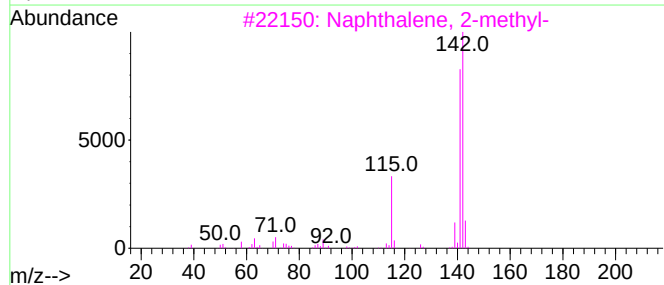
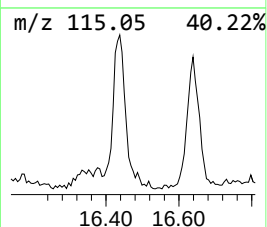
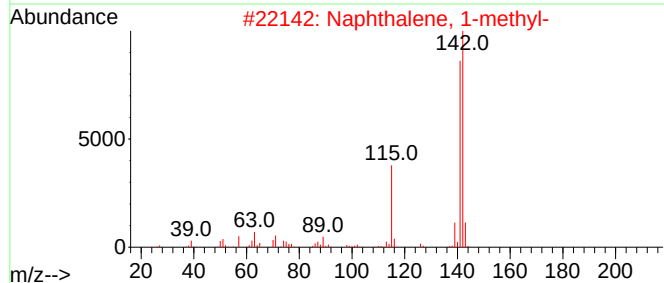
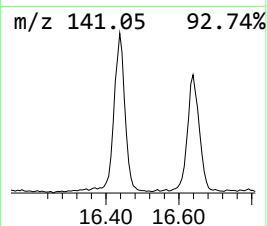
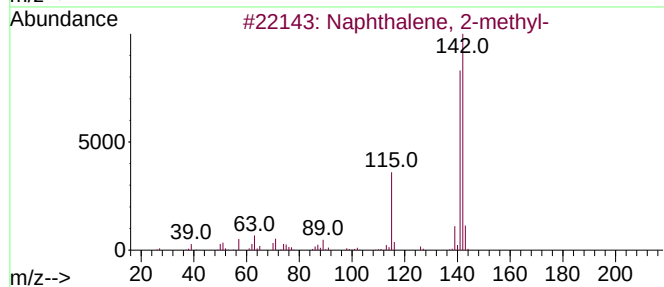
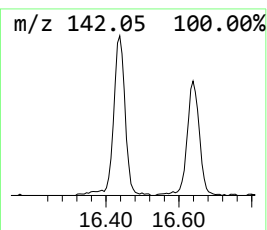
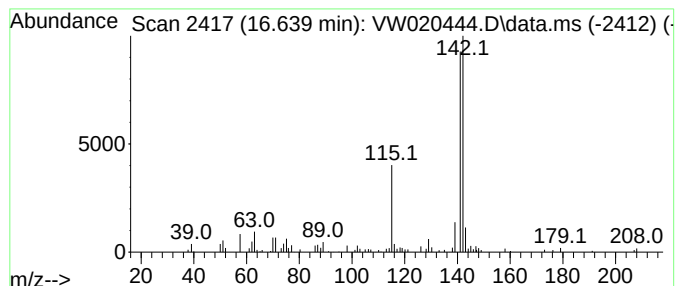
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM100821SMA.M
Quant Title : SFAM01.0

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

Peak Number 3 Naphthalene, 1-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.639	2.03 ug/L	73036	1,4-Dichlorobenzene-d4	13.560

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2-methyl-	142	C11H10	000091-57-6	94
2			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	94
3			Naphthalene, 2-methyl-	142	C11H10	000091-57-6	94
4			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	91
5			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	91



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TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Naphthalene, 2-...	16.438	2.5	ug/L	90389	3	13.560	900418	25.0
Naphthalene, 1-...	16.639	2.0	ug/L	73036	3	13.560	900418	25.0