LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM052220\
Data File : BM026101.D
Acq On : 23 May 2020 04:53
Operator : MA/SJ
Sample : L2737-13
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9B10

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % 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:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Title : SVOA CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	6.716	643	651	665	rVB	1989597	3666254	1.80%	0.358%
2	7.216	730	736	749	rVB	1769298	2724532	1.34%	0.266%
3	7.481	773	781	789	rBV3	1752484	3449801	1.70%	0.337%
4	7.804	828	836	840	rBV	2832982	4435284	2.18%	0.433%
5	7.863	840	846	853	rVV	7491653	11693114	5.75%	1.142%
6	7.957	853	862	867	rVV	9783609	15844992	7.79%	1.548%
7	8.022	867	873	884	rVB	2269708	3625078	1.78%	0.354%
8	8.281	910	917	922	rVV	8410065	15250074	7.50%	1.489%
9	8.340	922	927	939	rVB2	1917289	3838119	1.89%	0.375%
10	8.640	973	978	986	rVV2	936530	2164739	1.06%	0.211%
11	8.734	986	994	1001	rVB	2277370	3887557	1.91%	0.380%
12	8.910	1017	1024	1028	rBV	2129374	3672096	1.80%	0.359%
13	9.269	1080	1085	1093	rVB	869412	1642534	0.81%	0.160%
14	9.498	1113	1124	1128	rBV	8456537	23984093	11.79%	2.343%
15	9.687	1143	1156	1160	rBV	8966839	21267213	10.45%	2.077%
16	9.728	1160	1163	1171	rVV2	4665529	10530018	5.18%	1.028%
17	9.828	1171	1180	1188	rVV	13302269	27857655	13.69%	2.721%
18	9.922	1188	1196	1199	rVV2	1230770	2936017	1.44%	0.287%
19	9.963	1199	1203	1210	rVB	3181846	5385112	2.65%	0.526%
20	10.193	1231	1242	1250	rBV	3605699	6115284	3.01%	0.597%
21	10.387	1250	1275	1277	rBV4	40270757	203456864	100.00%	19.872%
22	10.463	1283	1288	1295	rVB	13745242	20190924	9.92%	1.972%
23	10.634	1309	1317	1324	rBV2	1294865	2685611	1.32%	0.262%
24	10.816	1337	1348	1352	rBV	2139038	3762220	1.85%	0.367%
25	11.116	1388	1399	1408	rBV2	4786503	10049810	4.94%	0.982%
26	11.292	1416	1429	1435	rBV3	1637217	3853668	1.89%	0.376%
27	11.357	1435	1440	1445	rBV	1902385	3301802	1.62%	0.322%
28	11.657	1483	1491	1500	rVB	3430731	6223262	3.06%	0.608%
29	11.769	1506	1510	1516	rVV2	617702	1431594	0.70%	0.140%
30	11.934	1527	1538	1544	rVV	8622735	14168778	6.96%	1.384%
31	12.039	1551	1556	1562	rVB2	2143399	3583718	1.76%	0.350%
32	12.157	1562	1576	1583	rBV	9086990	14904178	7.33%	1.456%
33	12.339	1602	1607	1612	rVB	2443403	3805795	1.87%	0.372%
34	12.875	1680	1698	1702	rBV7	431568	1534063	0.75%	0.150%

LSC Area Percent Report

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Operator : MA/SJ
Sample : L2737-13
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9B10

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % 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:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Title : SVOA CALIBRATION

35	12.934	1702	1708	1710	rBV2	1250928	1957105	0.96%	0.191%
36	12.963	1710	1713	1724	rVB2	1681161	3764180	1.85%	0.368%
37	13.286	1761	1768	1771	rBV4	1367201	2729625	1.34%	0.267%
38	13.457	1792	1797	1802	rVV	2320588	4293956	2.11%	0.419%
39	13.510	1802	1806	1811	rVV	1458120	2185560	1.07%	0.213%
40	13.563	1811	1815	1820	rVB	1993085	2728311	1.34%	0.266%
41	13.692	1832	1837	1840	rVV	1080316	1593840	0.78%	0.156%
42	13.822	1853	1859	1862	rBV	2414863	3578659	1.76%	0.350%
43	13.857	1862	1865	1874	rVB	1668477	2534552	1.25%	0.248%
44	14.139	1907	1913	1918	rBV3	2083903	3625726	1.78%	0.354%
45	14.216	1918	1926	1931	rVV	20795955	33127254	16.28%	3.236%
46	14.275	1931	1936	1941	rVV	2532773	3603214	1.77%	0.352%
47	14.345	1941	1948	1957	rVV	7369605	12007130	5.90%	1.173%
48	14.428	1957	1962	1971	rVB4	4184674	8455482	4.16%	0.826%
49	14.545	1971	1982	1992	rBV	9730576	17649991	8.68%	1.724%
50	14.645	1992	1999	2004	rVB	2430968	3648337	1.79%	0.356%
51	15.016	2038	2062	2065	rVV	11288901	50980134	25.06%	4.979%
52	15.092	2066	2075	2079	rVV	2069179	3512173	1.73%	0.343%
53	15.139	2079	2083	2088	rVV	3072712	4388869	2.16%	0.429%
54	15.198	2088	2093	2100	rVV	10948570	15735692	7.73%	1.537%
55	15.280	2100	2107	2111	rVV3	1190761	2325870	1.14%	0.227%
56	15.328	2111	2115	2116	rVV2	1059298	1490433	0.73%	0.146%
57	15.357	2116	2120	2125	rVV	3493167	5964877	2.93%	0.583%
58	15.445	2125	2135	2140	rVV4	2100127	5830458	2.87%	0.569%
59	15.545	2147	2152	2156	rVV	2351808	3274700	1.61%	0.320%
60	15.598	2156	2161	2164	rVV	1363522	2186637	1.07%	0.214%
61	15.639	2164	2168	2176	rVV2	1670648	2930805	1.44%	0.286%
62	15.875	2202	2208	2211	rBV	4628469	7416310	3.65%	0.724%
63	15.945	2211	2220	2224	rVB	5136119	14951379	7.35%	1.460%
64	16.110	2242	2248	2253	rBV2	3153531	6333055	3.11%	0.619%
65	16.186	2253	2261	2265	rBV3	5064821	10893811	5.35%	1.064%
66	16.369	2281	2292	2298	rVB3	8209622	26569572	13.06%	2.595%
67	16.704	2335	2349	2354	rBV2	8343793	22504011	11.06%	2.198%
68	16.833	2365	2371	2376	rBV2	1113866	2034469	1.00%	0.199%
69	16.892	2376	2381	2383	rBV2	2307255	3612062	1.78%	0.353%
70	16.939	2383	2389	2393	rVV	18368567	32168109	15.81%	3.142%
71	16.992	2393	2398	2401	rVV2	8243121	12818321	6.30%	1.252%

LSC Area Percent Report

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Sample : L2737-13
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9B10

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

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Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Title : SVOA CALIBRATION

72	17.022	2401	2403	2407	rVB	1659495	1886817	0.93%	0.184%
73	17.086	2407	2414	2416	rBV5	982322	1808028	0.89%	0.177%
74	17.174	2421	2429	2432	rBV	3277356	6361801	3.13%	0.621%
75	17.316	2432	2453	2459	rVB3	20029055	61985875	30.47%	6.054%
76	17.386	2459	2465	2466	rBV	2281971	3083639	1.52%	0.301%
77	17.421	2466	2471	2477	rVV	5399119	10863062	5.34%	1.061%
78	17.516	2478	2487	2495	rVB3	7505273	24984362	12.28%	2.440%
79	17.757	2524	2528	2534	rVB2	2933845	4638504	2.28%	0.453%
80	17.821	2534	2539	2542	rBV2	3380985	4982518	2.45%	0.487%
81	17.874	2542	2548	2552	rVB4	834068	1609457	0.79%	0.157%
82	17.927	2552	2557	2560	rBV2	2356471	4695390	2.31%	0.459%
83	18.121	2586	2590	2594	rVV3	1006755	1488173	0.73%	0.145%
84	18.257	2608	2613	2618	rVB3	1750424	2952547	1.45%	0.288%
85	18.310	2618	2622	2627	rVB2	1532205	2036779	1.00%	0.199%
86	18.957	2728	2732	2736	rVB	4417021	5598166	2.75%	0.547%
87	19.180	2765	2770	2775	rVV	6952092	9562604	4.70%	0.934%
88	19.292	2784	2789	2791	rBV2	3965840	5788922	2.85%	0.565%
89	19.315	2791	2793	2797	rVV	2657704	3122502	1.53%	0.305%
90	19.380	2797	2804	2808	rVB2	9309244	14021551	6.89%	1.370%
91	19.710	2856	2860	2864	rVB	1178809	1463700	0.72%	0.143%
92	19.810	2868	2877	2879	rBV	4659745	9519961	4.68%	0.930%
93	19.839	2879	2882	2886	rVB2	5118163	7198677	3.54%	0.703%
94	20.333	2962	2966	2975	rVB	1028854	1713730	0.84%	0.167%
95	20.468	2977	2989	2997	rVV3	1873535	5693133	2.80%	0.556%
96	21.086	3089	3094	3099	rVB2	2784294	3390735	1.67%	0.331%
97	21.574	3173	3177	3182	rVB	1316255	1701657	0.84%	0.166%
98	22.121	3264	3270	3280	rVB3	981956	1981248	0.97%	0.194%
99	23.074	3426	3432	3438	rBV	2607316	4309517	2.12%	0.421%
100	23.209	3449	3455	3467	rVB2	1729275	3071341	1.51%	0.300%

Sum of corrected areas: 1023840918

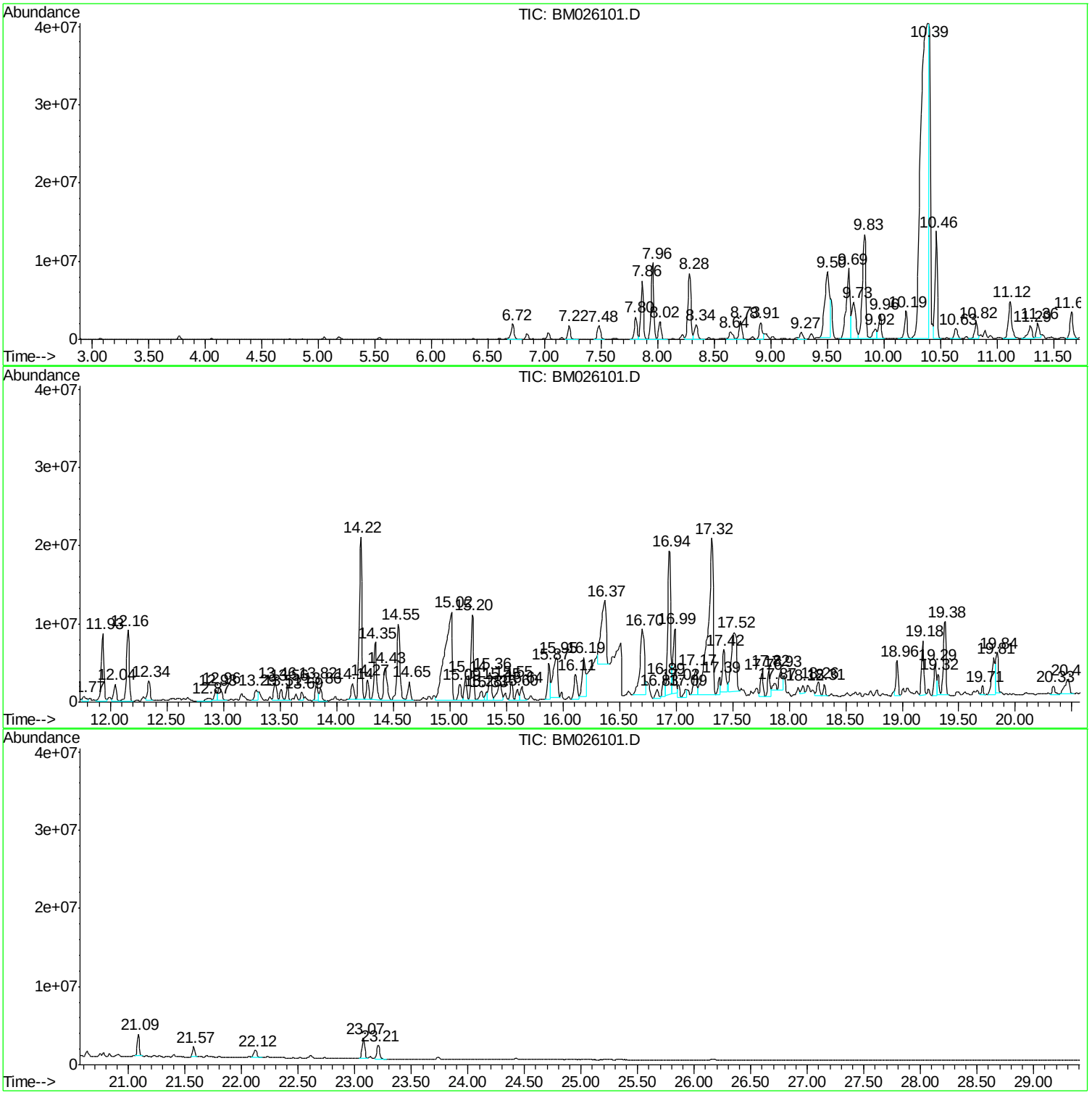
LSC Report - Integrated Chromatogram

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM052220\
Data File : BM026101.D
Acq On : 23 May 2020 04:53
Operator : MA/SJ
Sample : L2737-13
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9B10

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM052220\
Data File : BM026101.D
Acq On : 23 May 2020 04:53
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ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9B10

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

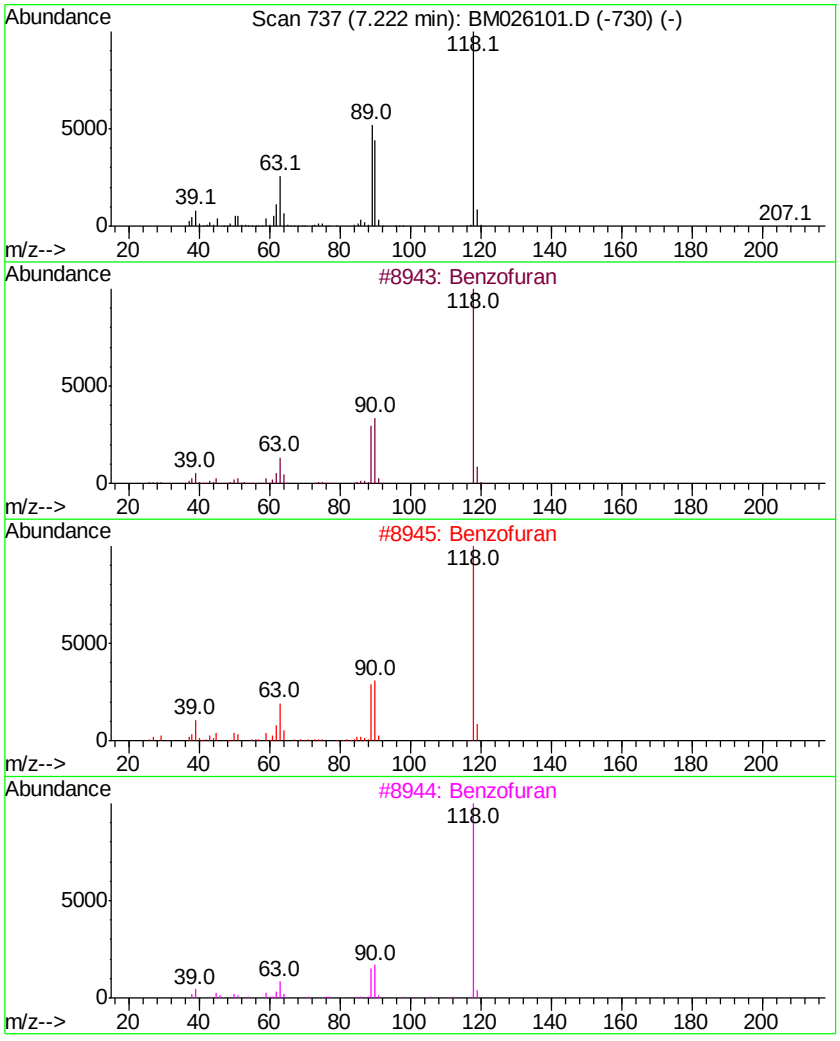
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

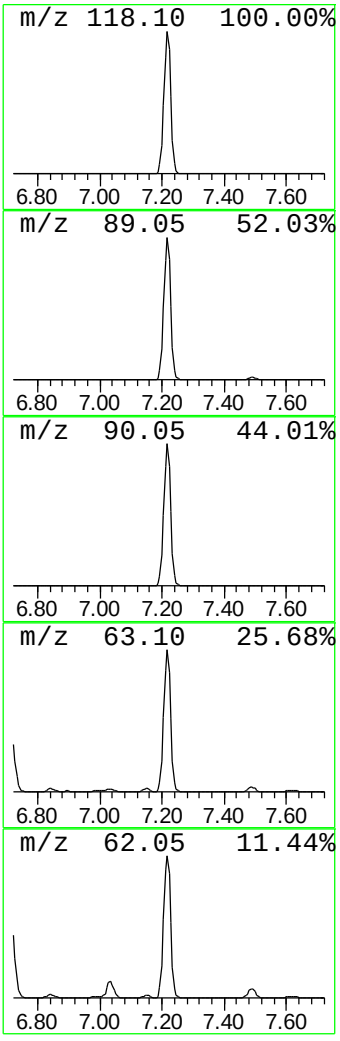
****

Peak Number 1 Benzofuran Concentration Rank 34

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.22	15.80 ng/ul	2724530	1,4-Dichlorobenzene-d4	7.49

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzofuran	118	C8H6O	000271-89-6	97
2	Benzofuran	118	C8H6O	000271-89-6	93
3	Benzofuran	118	C8H6O	000271-89-6	90
4	3-Hydroxyphenylacetylene	118	C8H6O	010401-11-3	74
5	2H-Cyclopenta[d]pyridazine	118	C7H6N2	000270-64-4	53





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM052220\
Data File : BM026101.D
Acq On : 23 May 2020 04:53
Operator : MA/SJ
Sample : L2737-13
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9B10

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

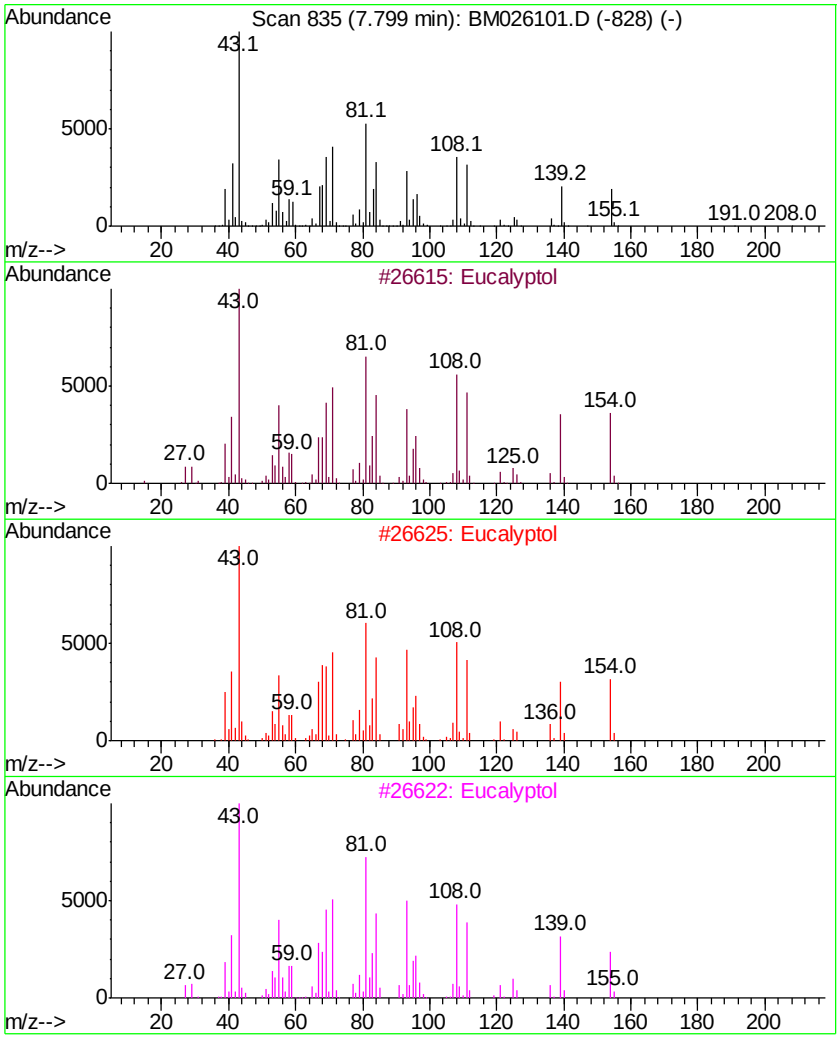
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

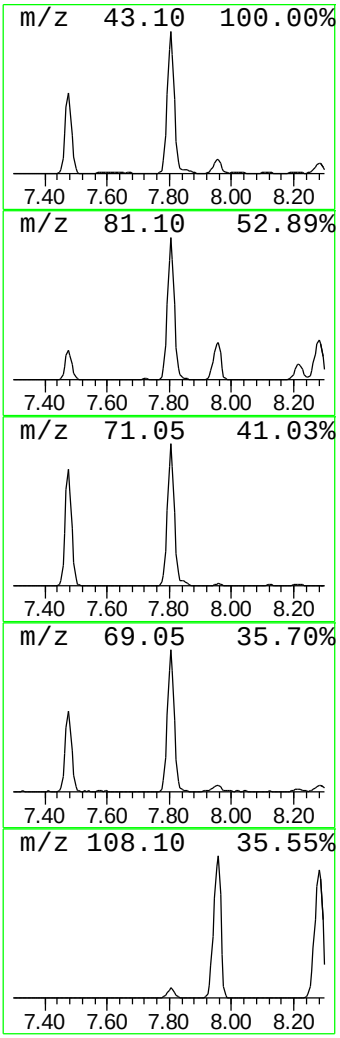
****

Peak Number 2 Eucalyptol Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.80	25.71 ng/ul	4435280	1,4-Dichlorobenzene-d4	7.49

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eucalyptol	154	C10H18O	000470-82-6	97
2	Eucalyptol	154	C10H18O	000470-82-6	94
3	Eucalyptol	154	C10H18O	000470-82-6	93
4	Eucalyptol	154	C10H18O	000470-82-6	91
5	Eucalyptol	154	C10H18O	000470-82-6	64





Library Search Compound Report

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Instrument : BNA_M
ClientSampleId : F9B10

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

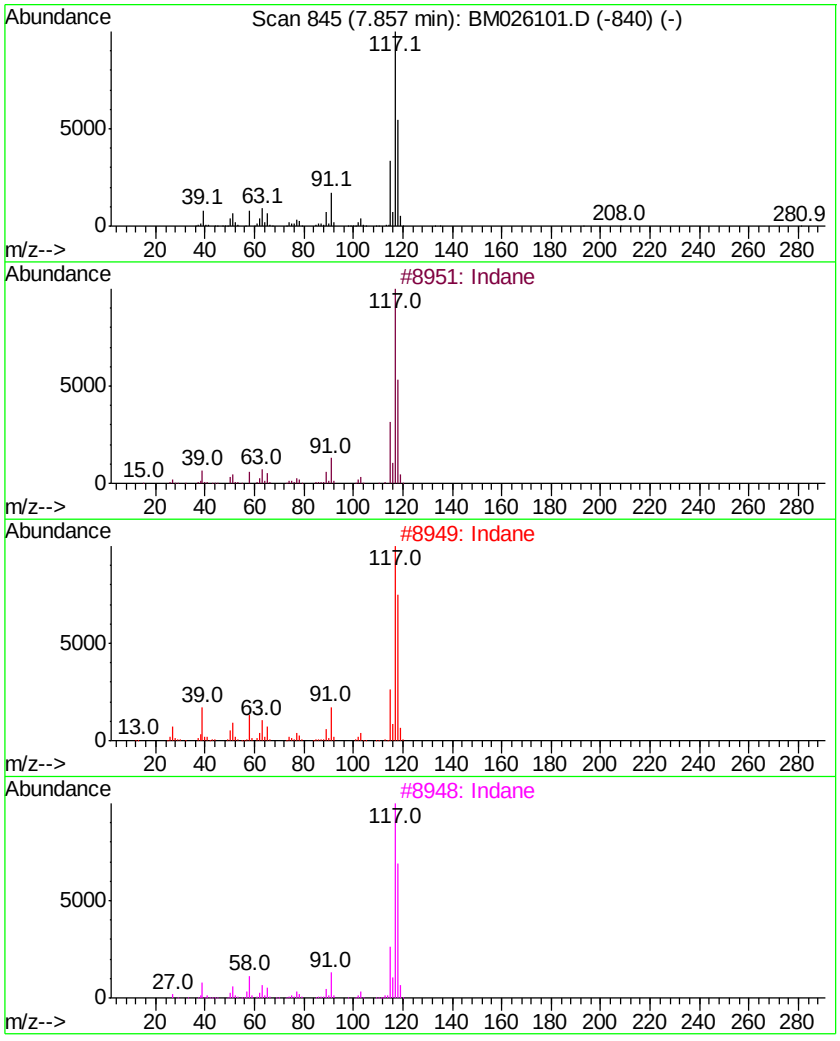
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

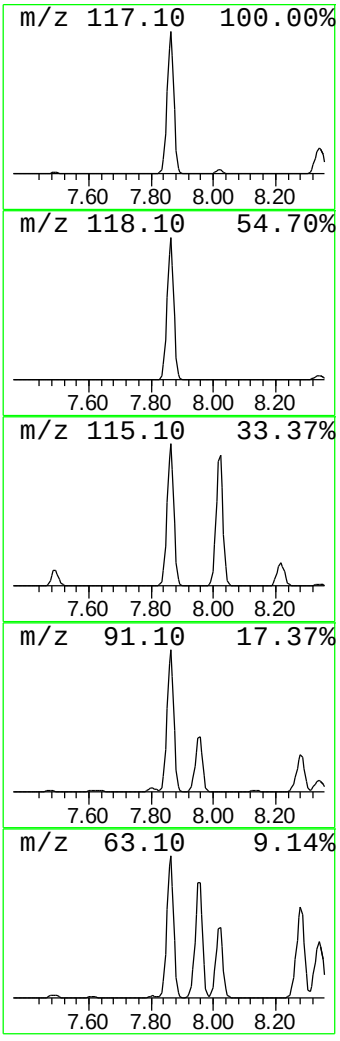
****

Peak Number 3 Benzene, 1-propenyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.86	67.79 ng/ul	11693100	1,4-Dichlorobenzene-d4	7.49

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	91
2	Indane	118	C9H10	000496-11-7	91
3	Indane	118	C9H10	000496-11-7	83
4	Benzene, 1-propenyl-	118	C9H10	000637-50-3	64
5	Benzene, 1-ethenyl-4-methyl-	118	C9H10	000622-97-9	64





Library Search Compound Report

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Data File : BM026101.D
Acq On : 23 May 2020 04:53
Operator : MA/SJ
Sample : L2737-13
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9B10

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

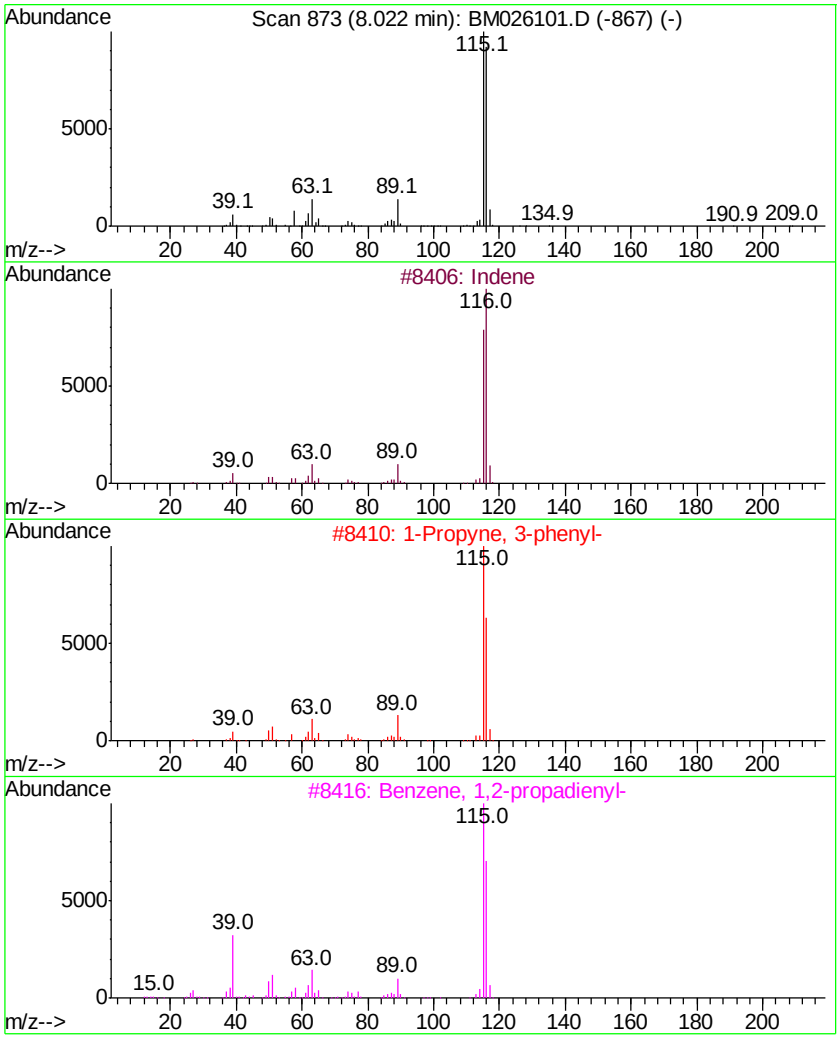
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

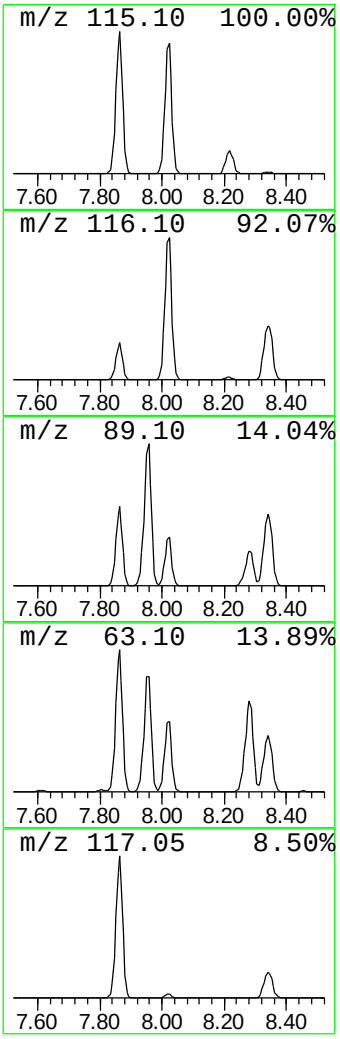
****

Peak Number 4 Indene Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.02	21.02 ng/ul	3625080	1,4-Dichlorobenzene-d4	7.49

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indene	116	C9H8	000095-13-6	97
2	1-Propyne, 3-phenyl-	116	C9H8	010147-11-2	94
3	Benzene, 1,2-propadienyl-	116	C9H8	002327-99-3	91
4	Benzene, 1-propynyl-	116	C9H8	000673-32-5	91
5	Benzene, 1-propynyl-	116	C9H8	000673-32-5	91





Library Search Compound Report

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ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9B10

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

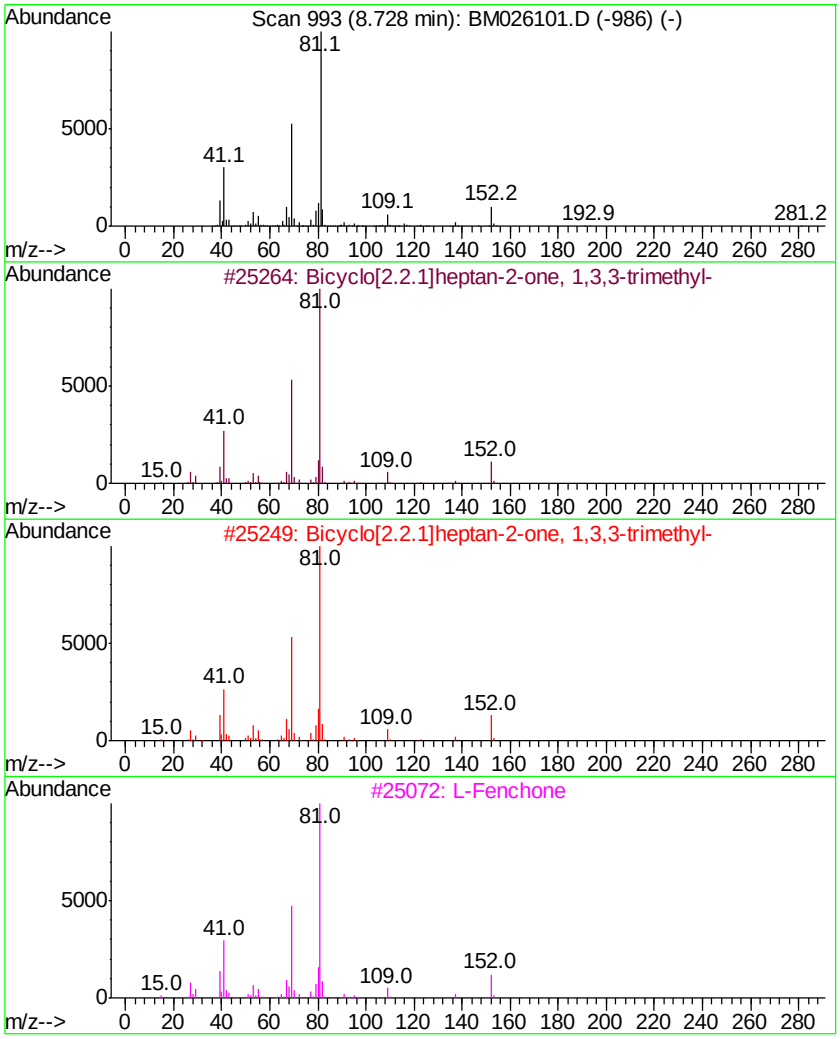
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

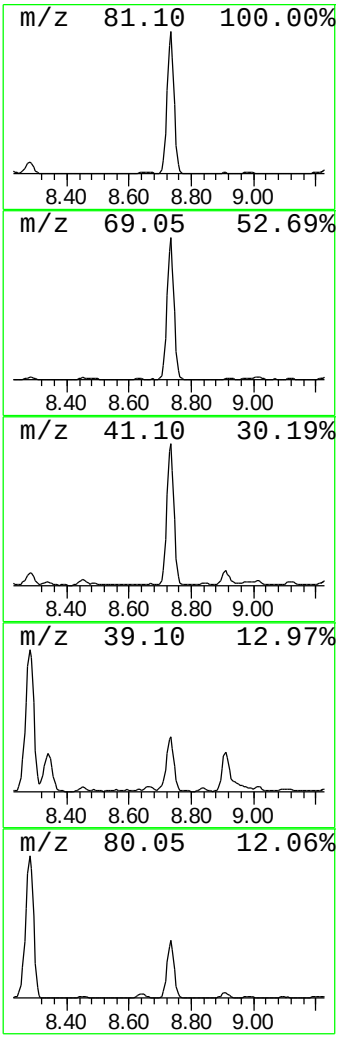
****

Peak Number 5 Bicyclo[2.2.1]heptan-2-one,... Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.73	22.54 ng/ul	3887560	1,4-Dichlorobenzene-d4	7.49

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Bicyclo[2.2.1]heptan-2-one, 1,3,...	152	C10H16O	001195-79-5	95
2	Bicyclo[2.2.1]heptan-2-one, 1,3,...	152	C10H16O	001195-79-5	95
3	L-Fenchone	152	C10H16O	007787-20-4	95
4	L-Fenchone	152	C10H16O	007787-20-4	94
5	Bicyclo[2.2.1]heptan-2-one, 1,3,...	152	C10H16O	001195-79-5	94





Library Search Compound Report

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Data File : BM026101.D
Acq On : 23 May 2020 04:53
Operator : MA/SJ
Sample : L2737-13
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9B10

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

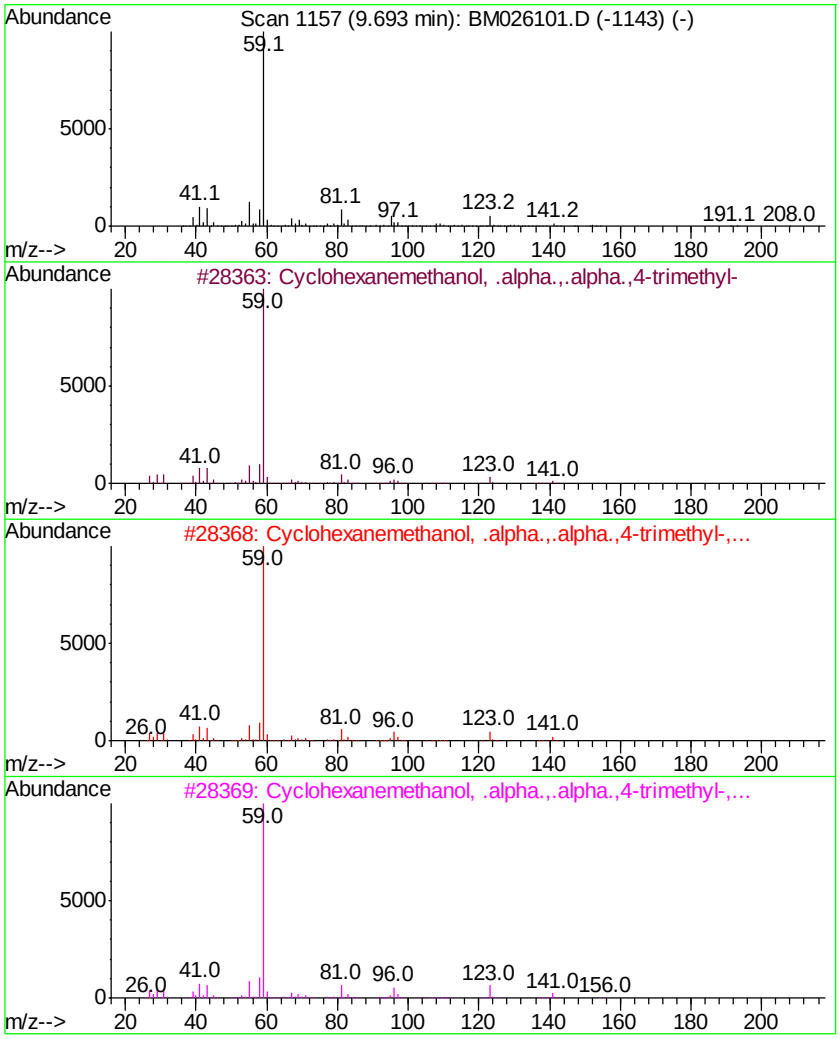
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

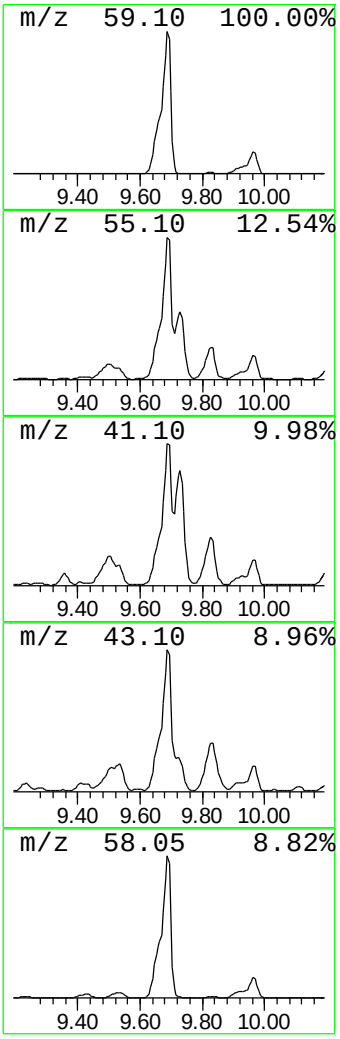
****

Peak Number 6 Cyclohexanemethanol, .alpha... Concentration Rank 50

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.69	2.09 ng/ul	21267200	Naphthalene-d8	10.29

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexanemethanol, .alpha.,.al...	156	C10H20O	000498-81-7	78
2	Cyclohexanemethanol, .alpha.,.al...	156	C10H20O	007322-63-6	78
3	Cyclohexanemethanol, .alpha.,.al...	156	C10H20O	005114-00-1	56
4	Cyclohexanemethanol, .alpha.,.al...	156	C10H20O	000498-81-7	47
5	2-Hydroxy-2,5-dimethyl-hept-6-en...	156	C9H16O2	1000192-55-8	45





Library Search Compound Report

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Sample : L2737-13
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9B10

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

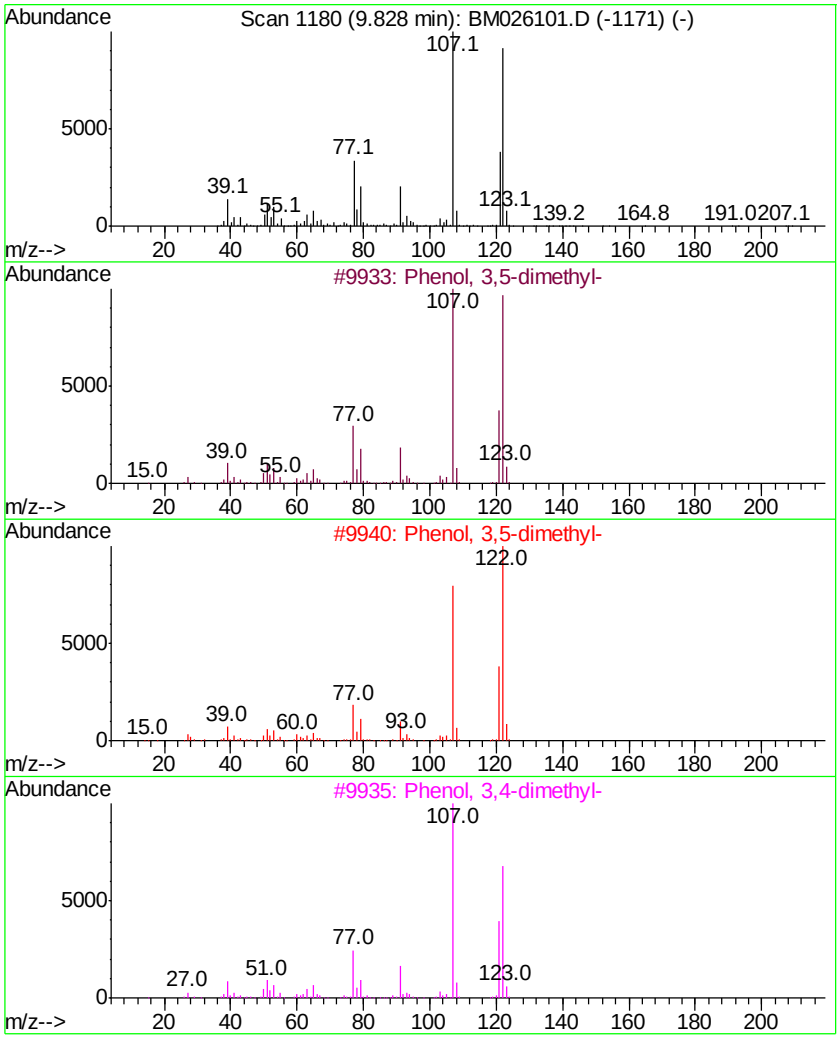
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

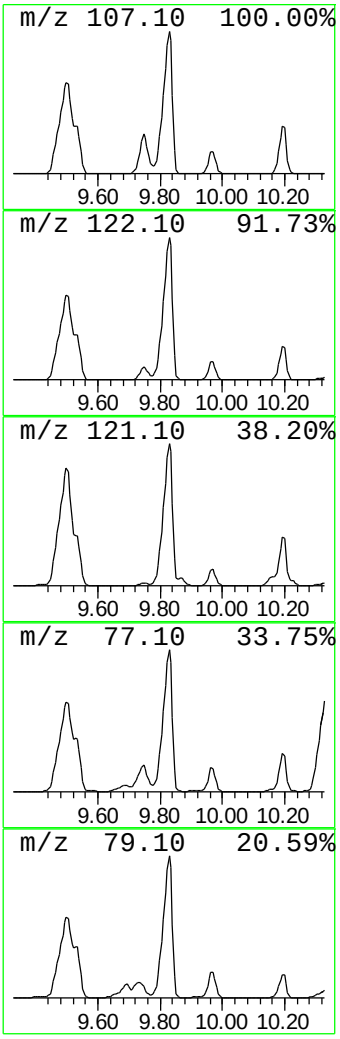
****

Peak Number 7 Phenol, 3,5-dimethyl- Concentration Rank 49

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.83	2.74 ng/ul	27857700	Naphthalene-d8	10.29

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 3,5-dimethyl-	122	C8H10O	000108-68-9	96
2	Phenol, 3,5-dimethyl-	122	C8H10O	000108-68-9	95
3	Phenol, 3,4-dimethyl-	122	C8H10O	000095-65-8	94
4	Phenol, 2,5-dimethyl-	122	C8H10O	000095-87-4	93
5	Phenol, 2,4-dimethyl-	122	C8H10O	000105-67-9	93





Library Search Compound Report

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Acq On : 23 May 2020 04:53
Operator : MA/SJ
Sample : L2737-13
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9B10

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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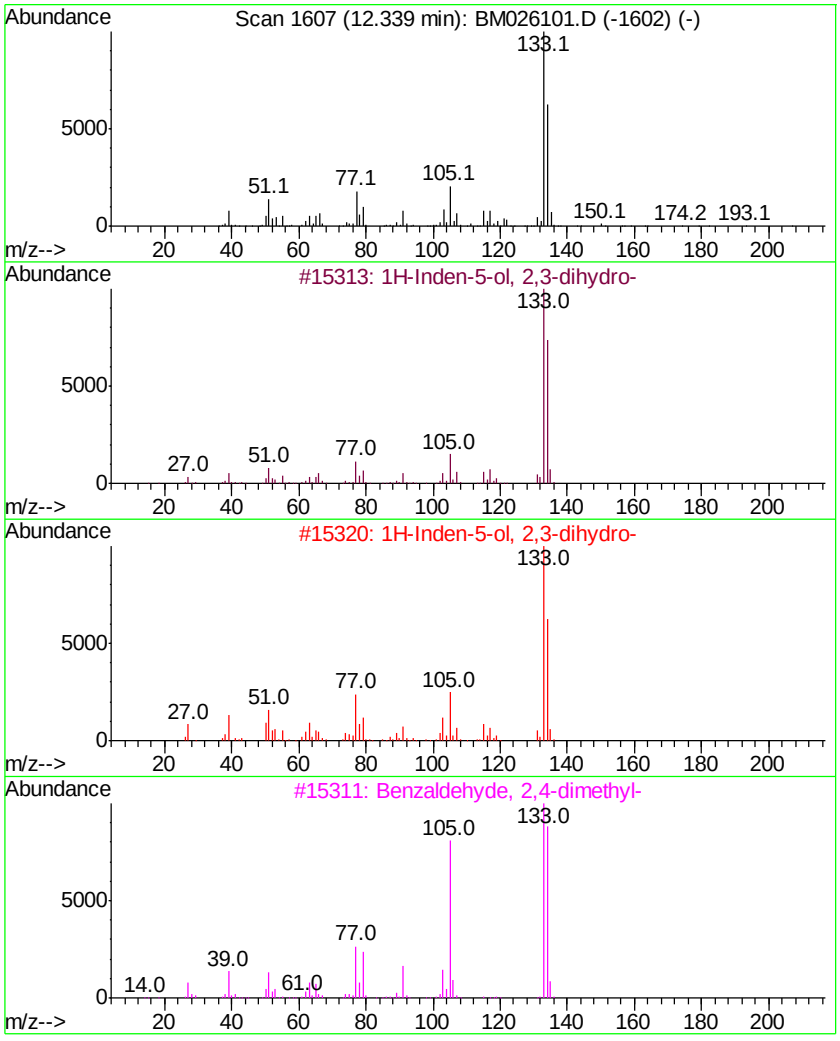
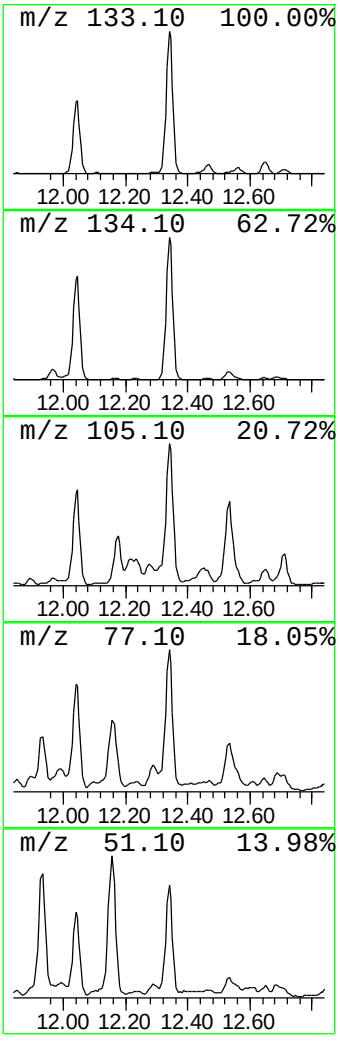
Peak Number 8 1H-Inden-5-ol, 2,3-dihydro- Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.34	20.99 ng/ul	3805800	Acenaphthene-d10	14.14

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Inden-5-ol, 2,3-dihydro-	134	C9H10O	001470-94-6	94
2	1H-Inden-5-ol, 2,3-dihydro-	134	C9H10O	001470-94-6	94
3	Benzaldehyde, 2,4-dimethyl-	134	C9H10O	015764-16-6	80
4	Benzaldehyde, 2,4-dimethyl-	134	C9H10O	015764-16-6	72
5	Benzaldehyde, 3,4-dimethyl-	134	C9H10O	005973-71-7	59





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Acq On : 23 May 2020 04:53
Operator : MA/SJ
Sample : L2737-13
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9B10

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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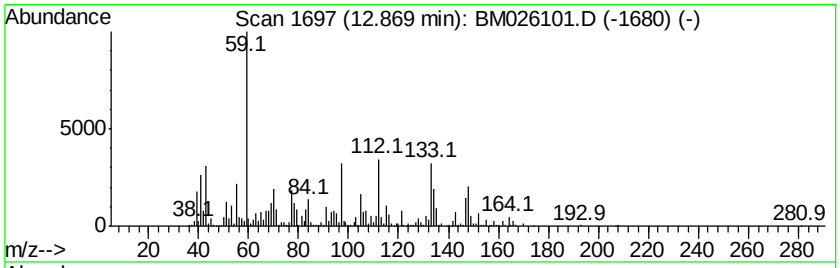
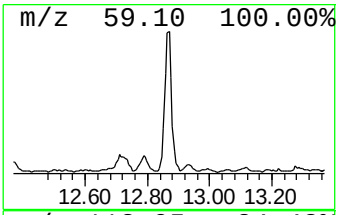
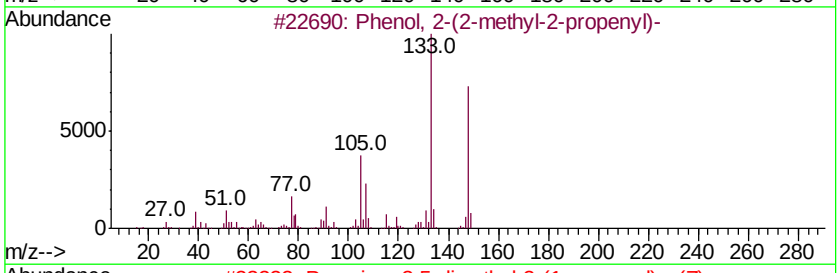
Peak Number 9 unknown-01 Concentration Rank 47

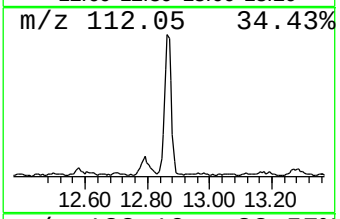
R.T.	EstConc	Area	Relative to ISTD	R.T.
12.87	8.46 ng/ul	1534060	Acenaphthene-d10	14.14

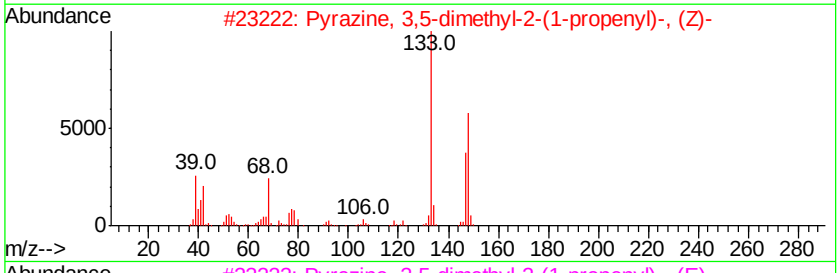
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 2-(2-methyl-2-propenyl)-	148	C10H12O	020944-88-1	38
2	Pyrazine, 3,5-dimethyl-2-(1-prop...	148	C9H12N2	055138-74-4	30
3	Pyrazine, 3,5-dimethyl-2-(1-prop...	148	C9H12N2	055138-78-8	30
4	Benzene, 1-ethyl-3-(1-methylethyl)-	148	C11H16	004920-99-4	25
5	3-(4-Pyridyl)acrylaldehyde	133	C8H7NO	026505-36-2	25

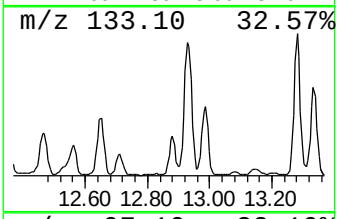


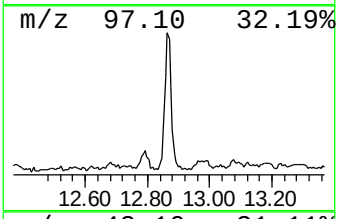


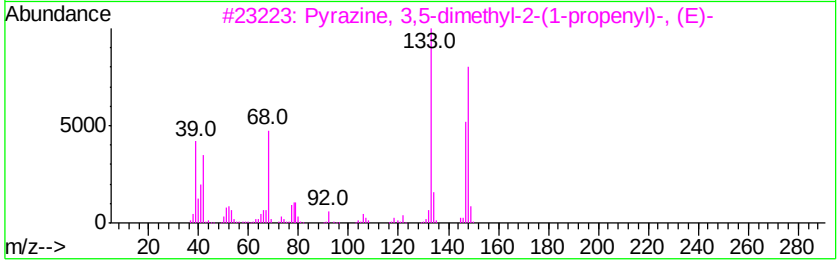


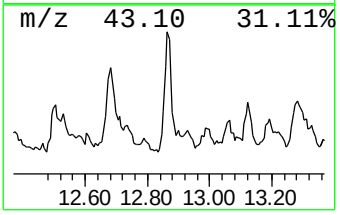












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Data File : BM026101.D
Acq On : 23 May 2020 04:53
Operator : MA/SJ
Sample : L2737-13
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9B10

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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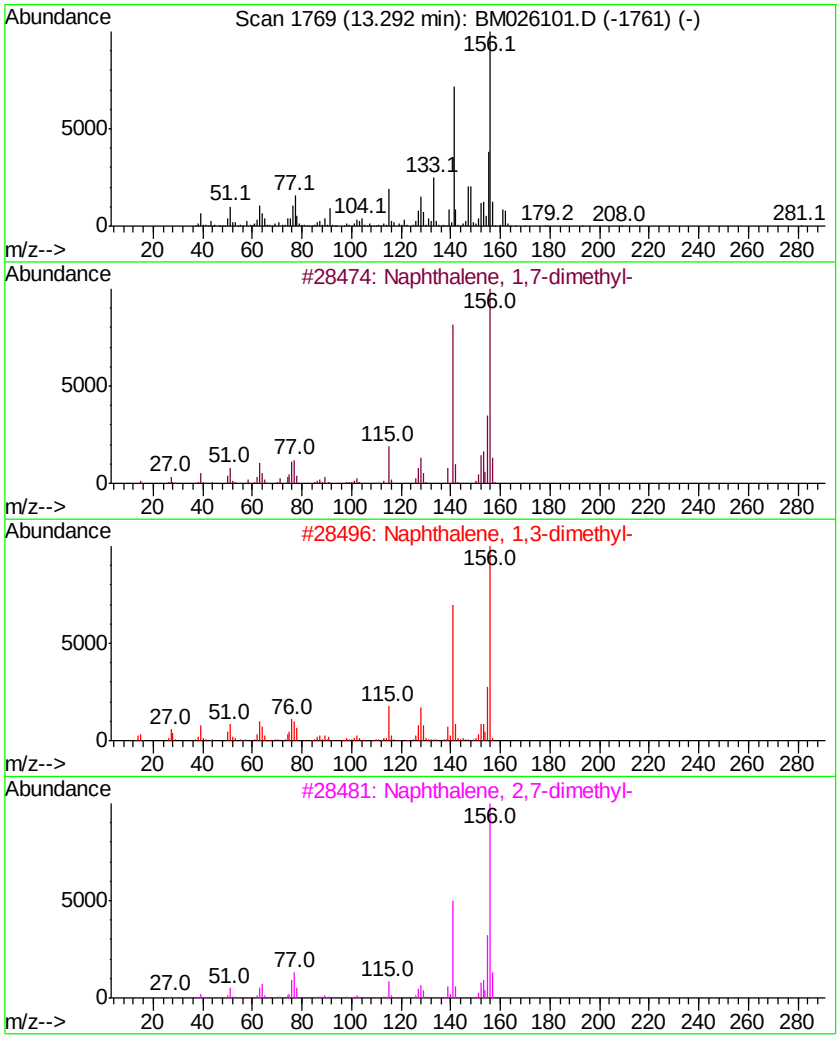
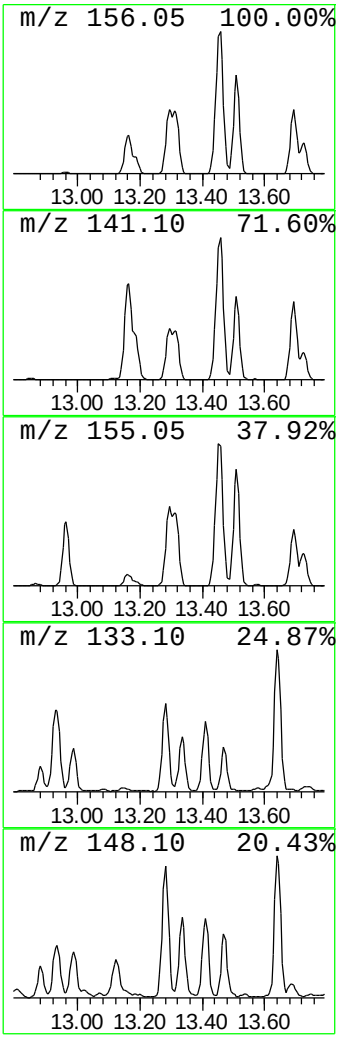
Peak Number 10 Naphthalene, 2,7-dimethyl- Concentration Rank 35

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.29	15.06 ng/ul	2729630	Acenaphthene-d10	14.14

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	95
2	Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	95
3	Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	95
4	Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	95
5	Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	95





Library Search Compound Report

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Acq On : 23 May 2020 04:53
Operator : MA/SJ
Sample : L2737-13
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9B10

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

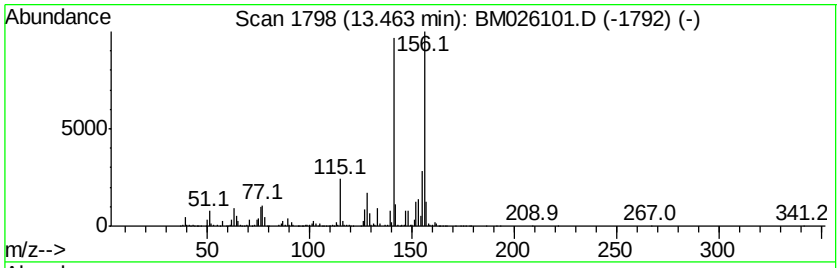
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

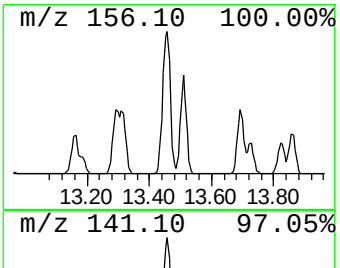
****

Peak Number 11 Naphthalene, 2,3-dimethyl- Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.46	23.69 ng/ul	4293960	Acenaphthene-d10	14.14

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	98
2	Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	98
3	Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	98
4	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
5	Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	97





Library Search Compound Report

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Sample : L2737-13
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9B10

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

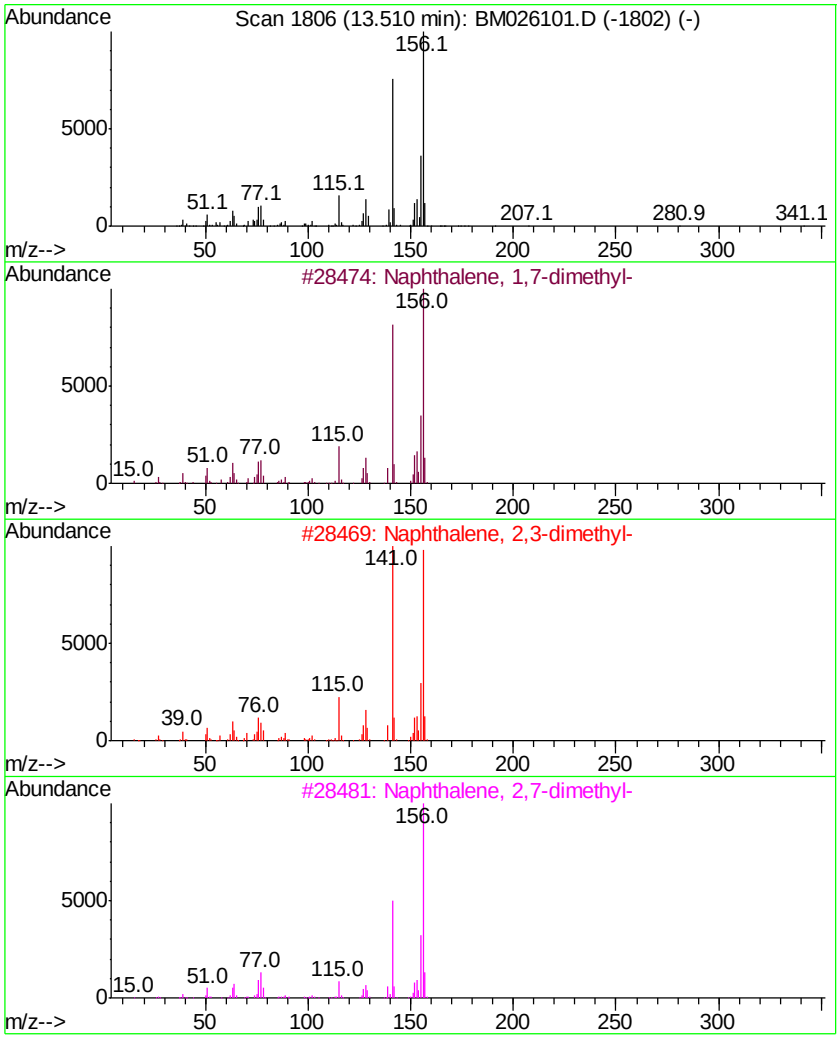
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

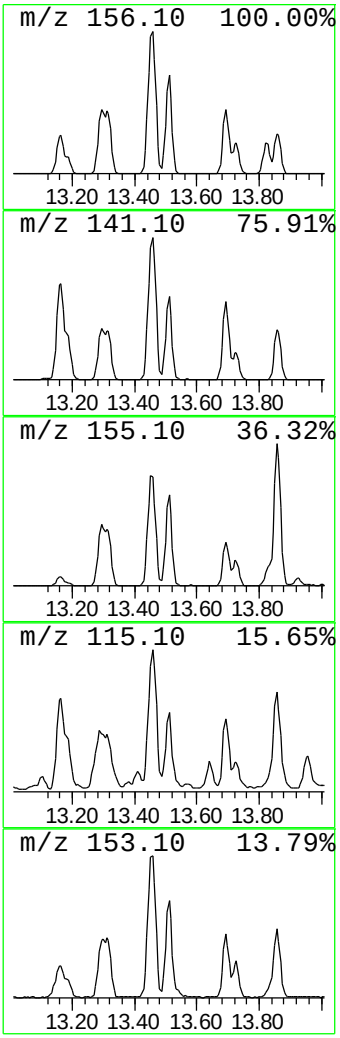
****

Peak Number 12 Naphthalene, 1,7-dimethyl- Concentration Rank 37

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.51	12.06 ng/ul	2185560	Acenaphthene-d10	14.14

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	98
2	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
3	Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	97
4	Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	97
5	Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	97





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Sample : L2737-13
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9B10

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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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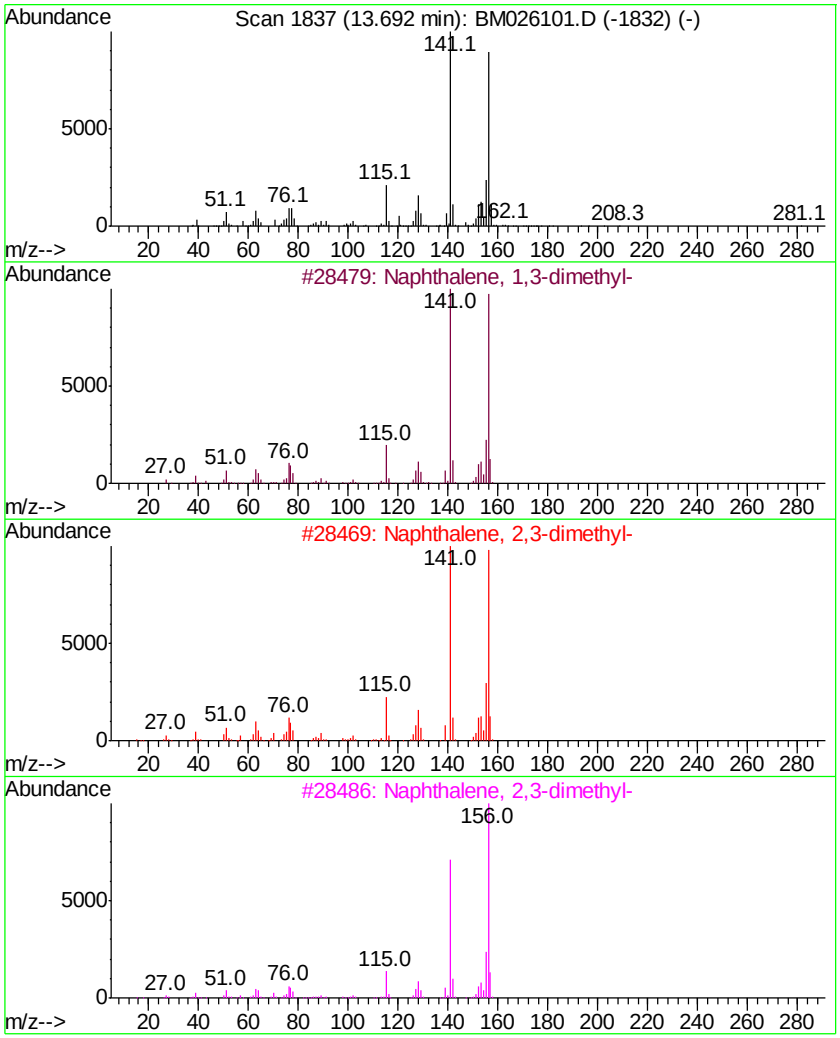
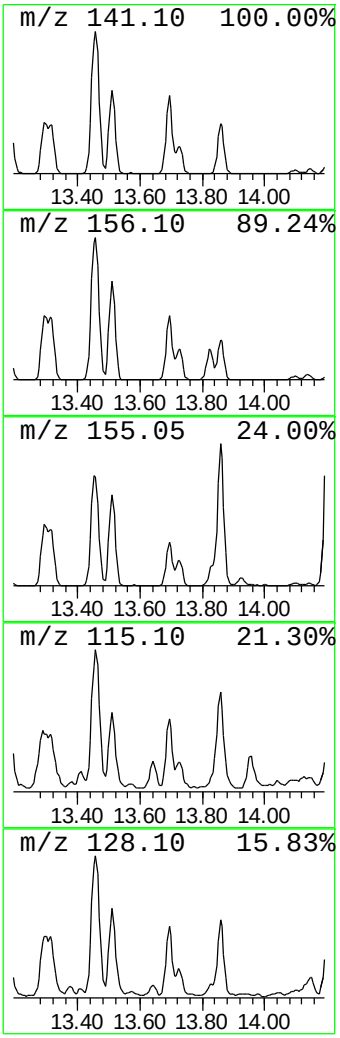
Peak Number 13 Naphthalene, 1,3-dimethyl- Concentration Rank 45

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.69	8.79 ng/ul	1593840	Acenaphthene-d10	14.14

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	98
2	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	98
3	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
4	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
5	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97





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ClientSampleId : F9B10

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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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Peak Number 14 2-Naphthalenecarbonitrile Concentration Rank 29

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.27	19.88 ng/ul	3603210	Acenaphthene-d10	14.14

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Naphthalenecarbonitrile	153	C11H7N	000613-46-7	97
2	1-Naphthalenecarbonitrile	153	C11H7N	000086-53-3	95
3	1-Naphthalenecarbonitrile	153	C11H7N	000086-53-3	94
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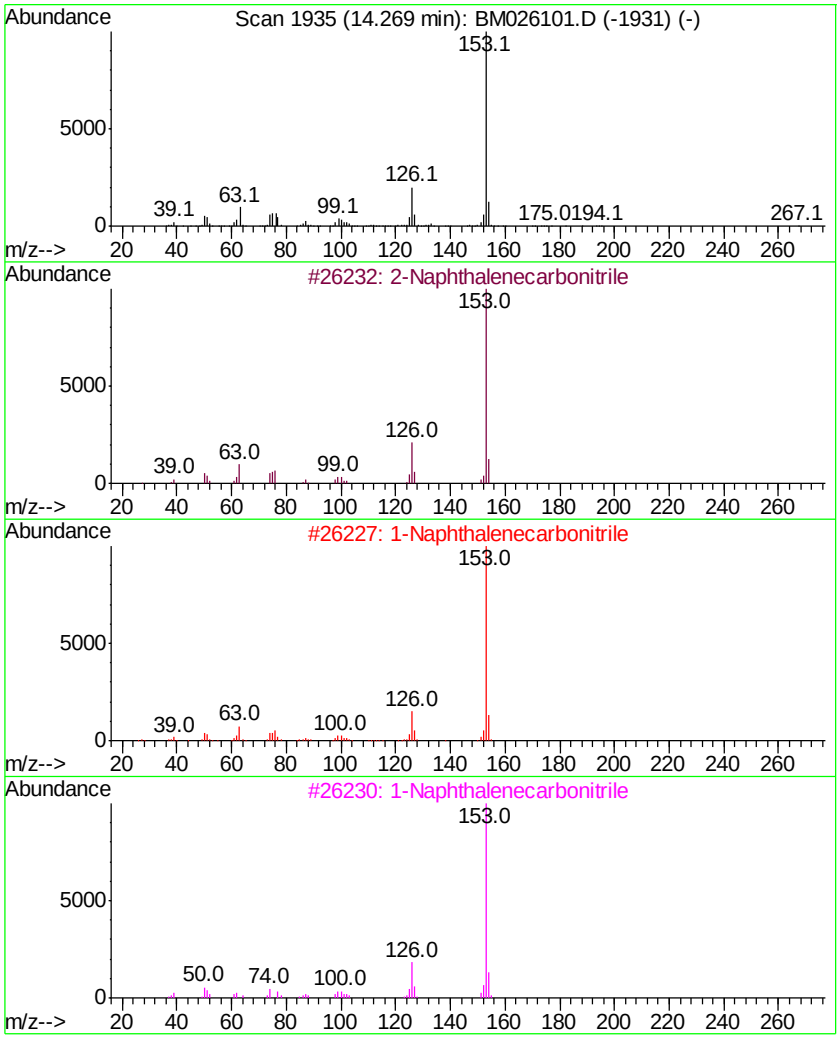

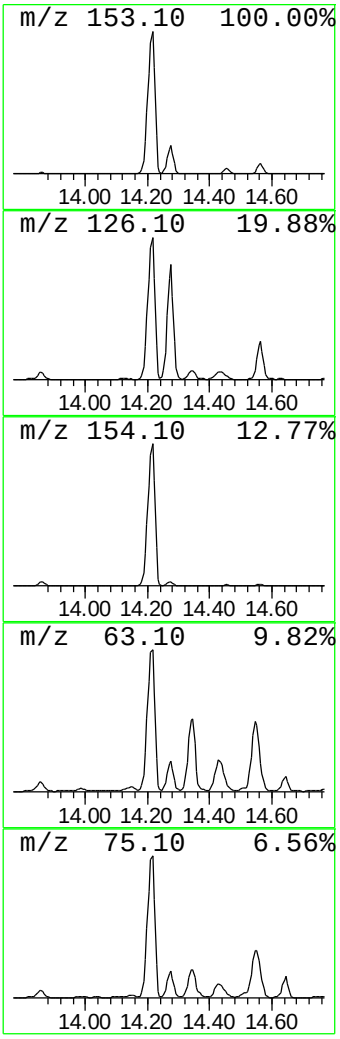

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Instrument : BNA_M
ClientSampleId : F9B10

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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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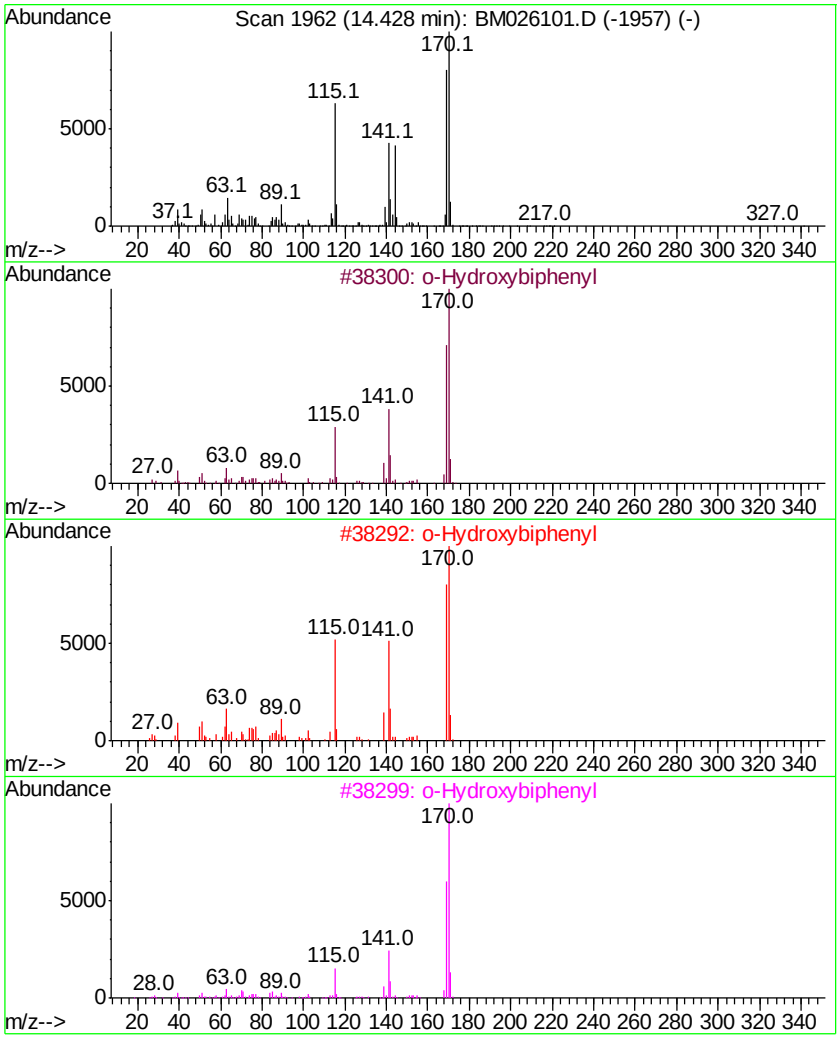
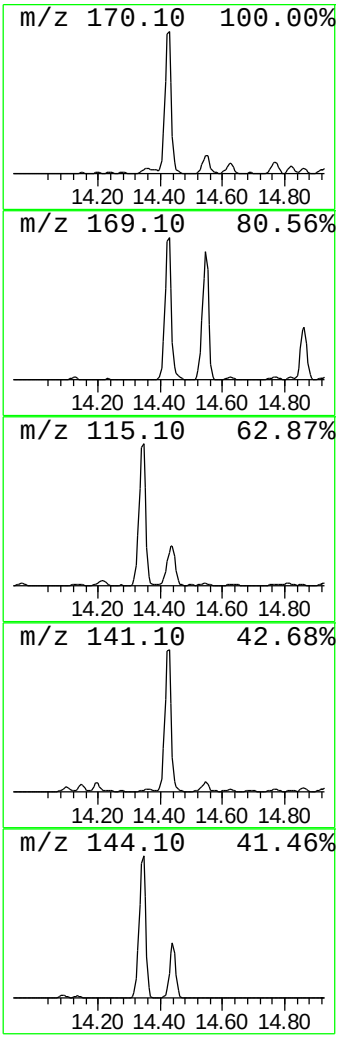
Peak Number 15 o-Hydroxybiphenyl Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.43	46.64 ng/ul	8455480	Acenaphthene-d10	14.14

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	o-Hydroxybiphenyl	170	C12H10O	000090-43-7	93
2	o-Hydroxybiphenyl	170	C12H10O	000090-43-7	76
3	o-Hydroxybiphenyl	170	C12H10O	000090-43-7	70
4	o-Hydroxybiphenyl	170	C12H10O	000090-43-7	70
5	o-Hydroxybiphenyl	170	C12H10O	000090-43-7	68





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Acq On : 23 May 2020 04:53
Operator : MA/SJ
Sample : L2737-13
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9B10

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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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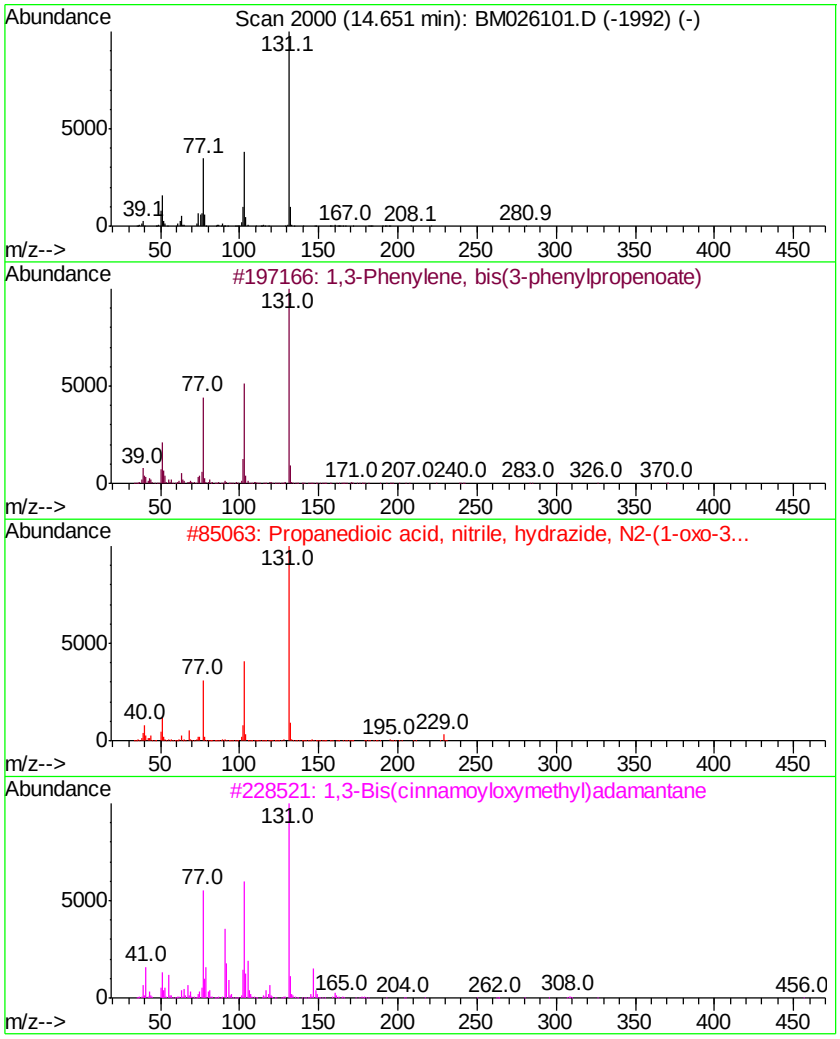
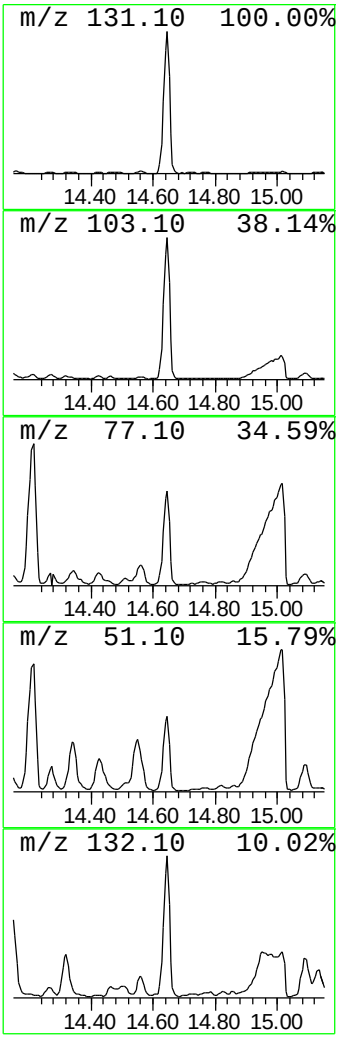
Peak Number 16 1,3-Phenylene, bis(3-phenyl... Concentration Rank 28

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.65	20.12 ng/ul	3648340	Acenaphthene-d10	14.14

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,3-Phenylene, bis(3-phenylprope...	370	C24H18O4	1000259-62-4	83
2	Propanedioic acid, nitrile, hydr...	229	C12H11N3O2	294878-35-6	83
3	1,3-Bis(cinnamoyloxymethyl)adama...	456	C30H32O4	303797-58-2	78
4	Cinnamaldehyde, (E)-	132	C9H8O	014371-10-9	78
5	Vinyl trans-cinnamate	174	C11H10O2	017719-70-9	74





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Sample : L2737-13
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9B10

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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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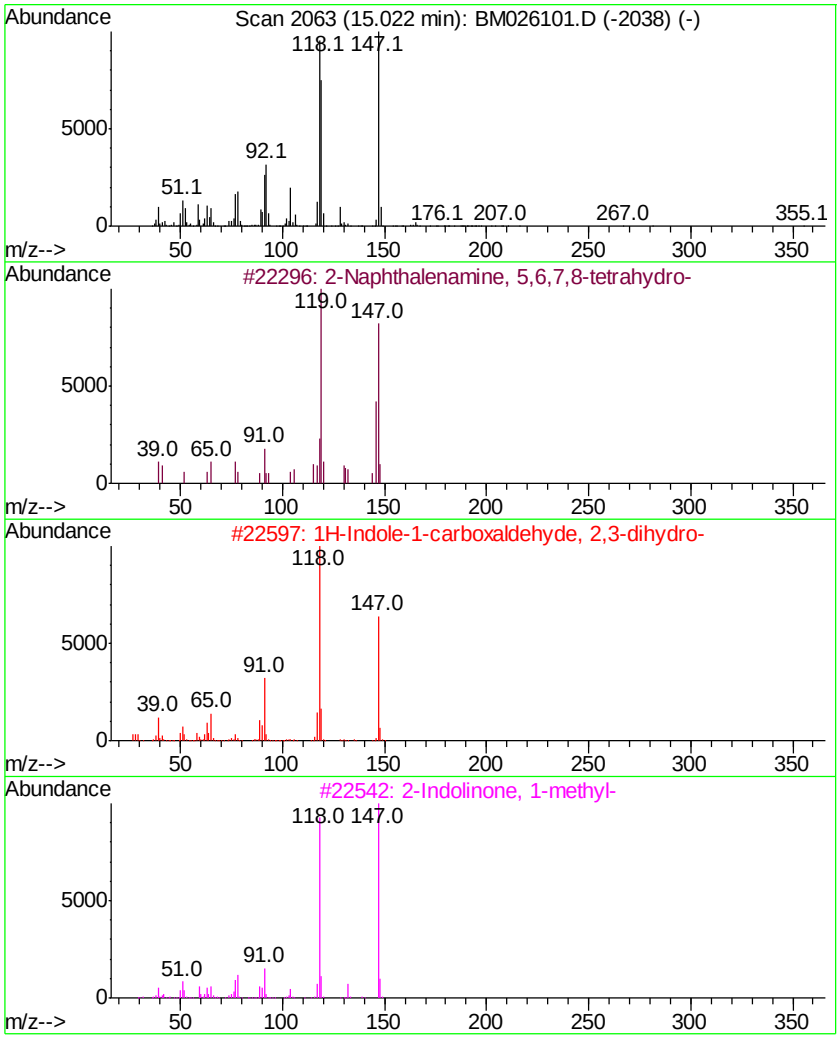
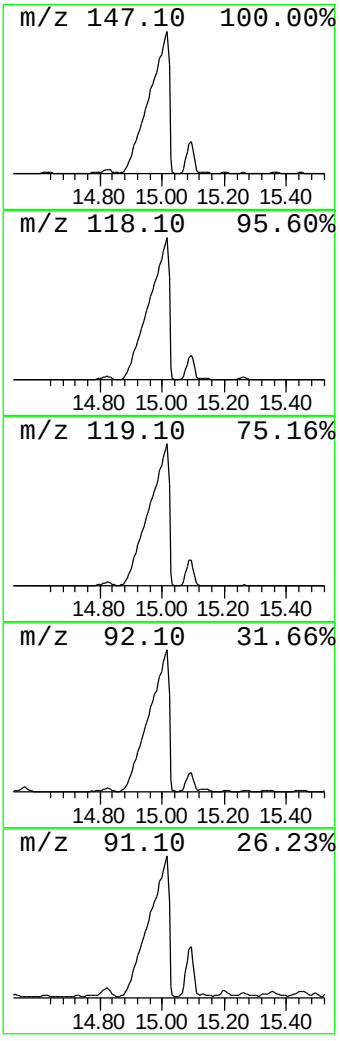
Peak Number 17 2-Naphthalenamine, 5,6,7,8-... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.02	281.21 ng/ul	50980100	Acenaphthene-d10	14.14

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Naphthalenamine, 5,6,7,8-tetra...	147	C10H13N	002217-43-8	64
2	1H-Indole-1-carboxaldehyde, 2,3-...	147	C9H9NO	002861-59-8	60
3	2-Indolinone, 1-methyl-	147	C9H9NO	000061-70-1	58
4	1H-Indole-1-carboxaldehyde, 2,3-...	147	C9H9NO	002861-59-8	55
5	(2-Benzimidazolyl)methylamine	147	C8H9N3	005805-57-2	52





Library Search Compound Report

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ClientSampleId : F9B10

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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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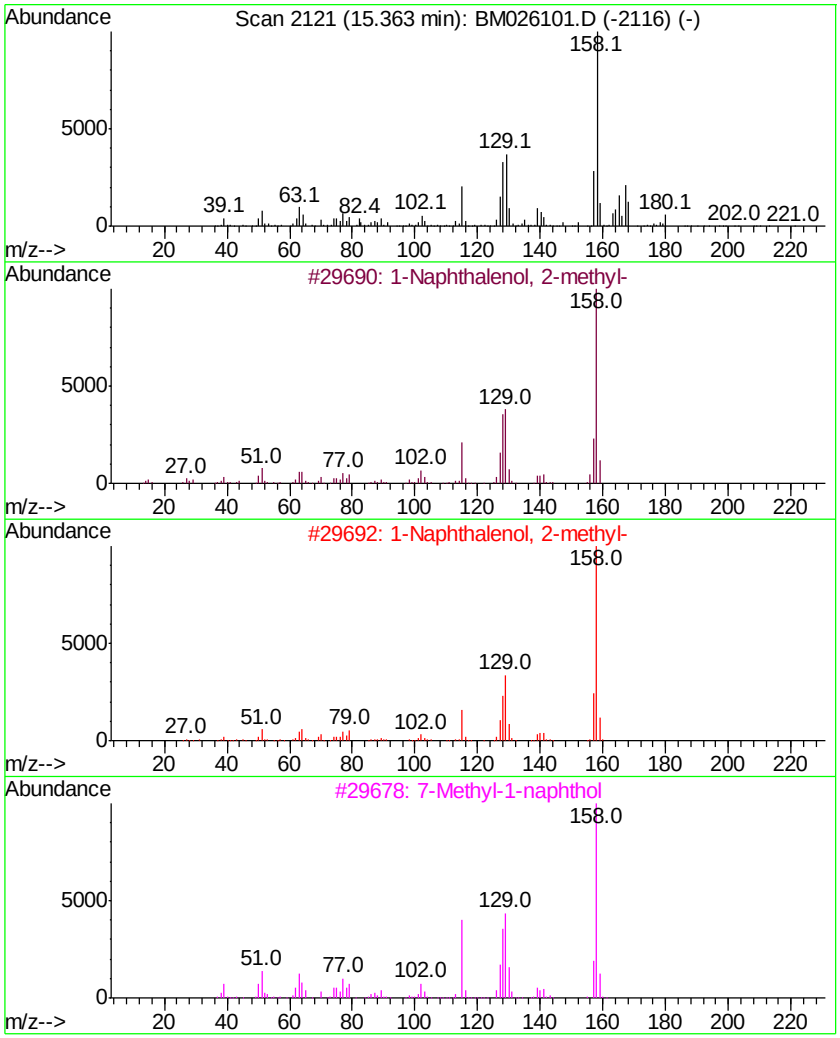
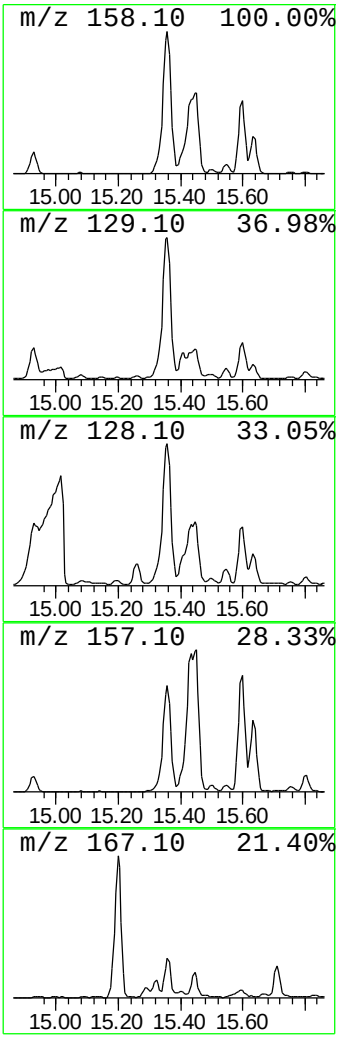
Peak Number 18 1-Naphthalenol, 2-methyl- Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.36	32.90 ng/ul	5964880	Acenaphthene-d10	14.14

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Naphthalenol, 2-methyl-	158	C11H10O	007469-77-4	96
2	1-Naphthalenol, 2-methyl-	158	C11H10O	007469-77-4	94
3	7-Methyl-1-naphthol	158	C11H10O	006939-33-9	93
4	Cinnoline, 3,4-dimethyl-	158	C10H10N2	003929-83-7	62
5	1-Naphthalenol, 4-methyl-	158	C11H10O	010240-08-1	47





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Sample : L2737-13
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9B10

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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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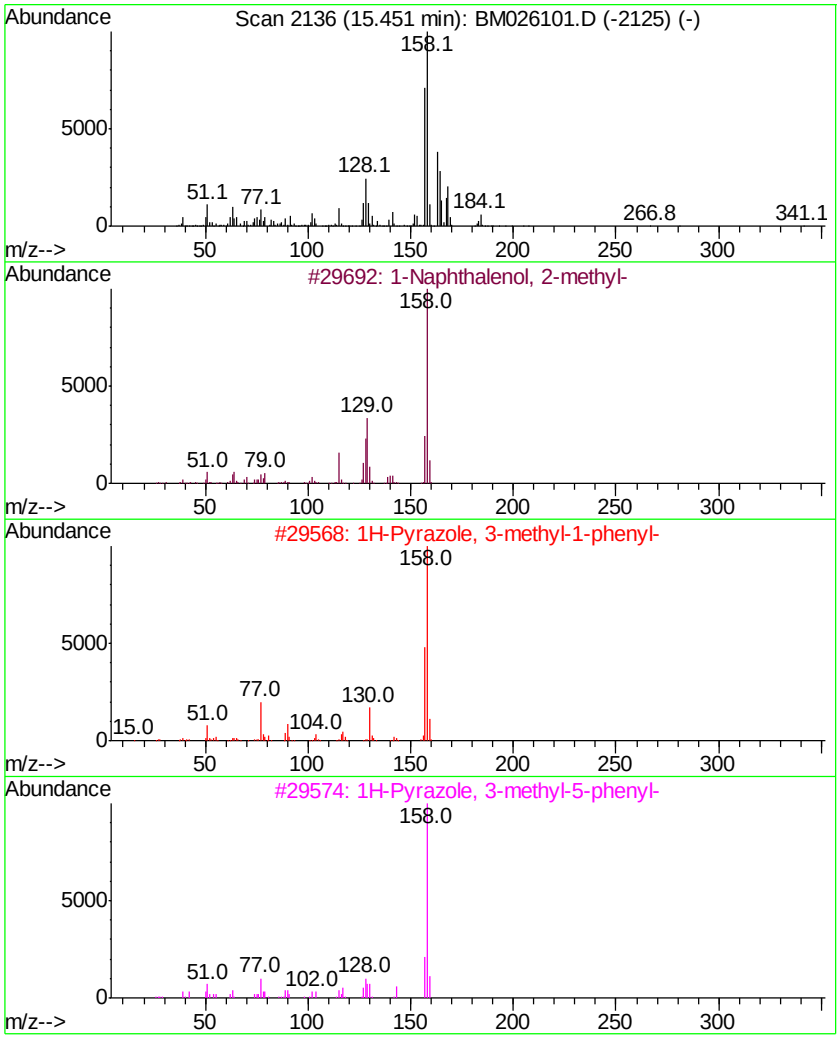
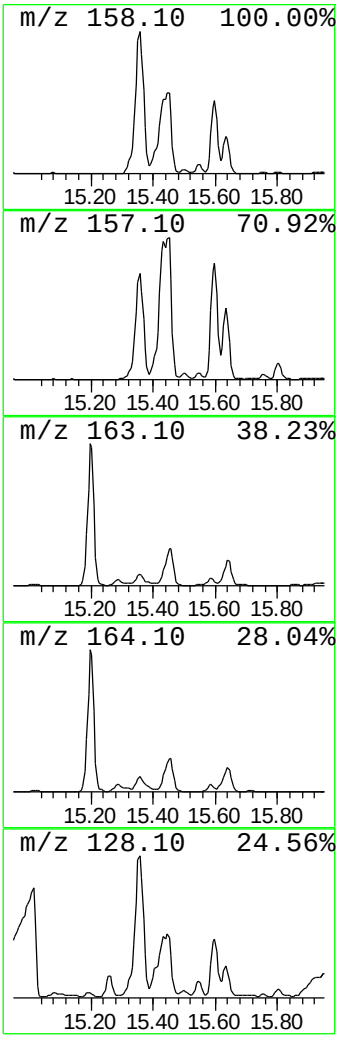
Peak Number 19 unknown-02 Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.45	32.16 ng/ul	5830460	Acenaphthene-d10	14.14

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Naphthalenol, 2-methyl-	158	C11H10O	007469-77-4	43
2	1H-Pyrazole, 3-methyl-1-phenyl-	158	C10H10N2	001128-54-7	38
3	1H-Pyrazole, 3-methyl-5-phenyl-	158	C10H10N2	003347-62-4	38
4	Pyrazole, 4-methyl-3-phenyl-	158	C10H10N2	013808-62-3	38
5	Benzene, 1,3-bis(1-methylethenyl)-	158	C12H14	003748-13-8	38





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Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9B10

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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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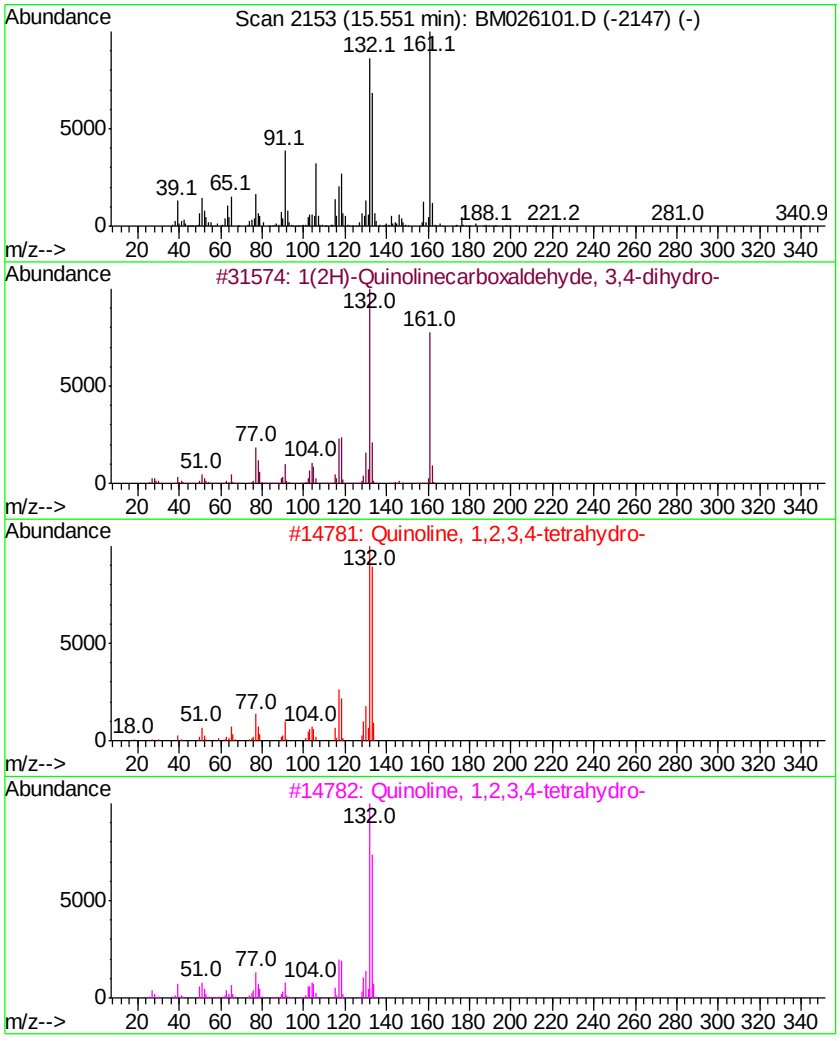
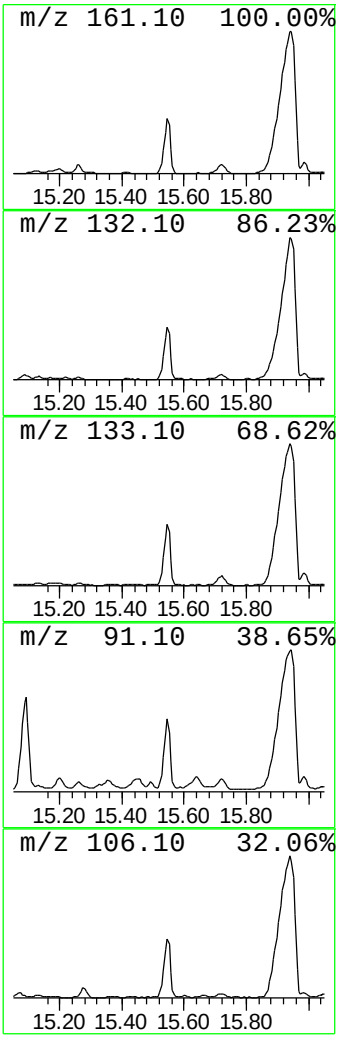
Peak Number 20 1(2H)-Quinolinecarboxaldehy... Concentration Rank 30

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.55	18.13 ng/ul	3274700	Phenanthrene-d10	16.89

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1(2H)-Quinolinecarboxaldehyde, 3...	161	C10H11NO	002739-16-4	68
2	Quinoline, 1,2,3,4-tetrahydro-	133	C9H11N	000635-46-1	55
3	Quinoline, 1,2,3,4-tetrahydro-	133	C9H11N	000635-46-1	55
4	5,6,7,8-Tetrahydroisoquinoline	133	C9H11N	1000379-32-5	53
5	Quinoline, 1,2,3,4-tetrahydro-	133	C9H11N	000635-46-1	49





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Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9B10

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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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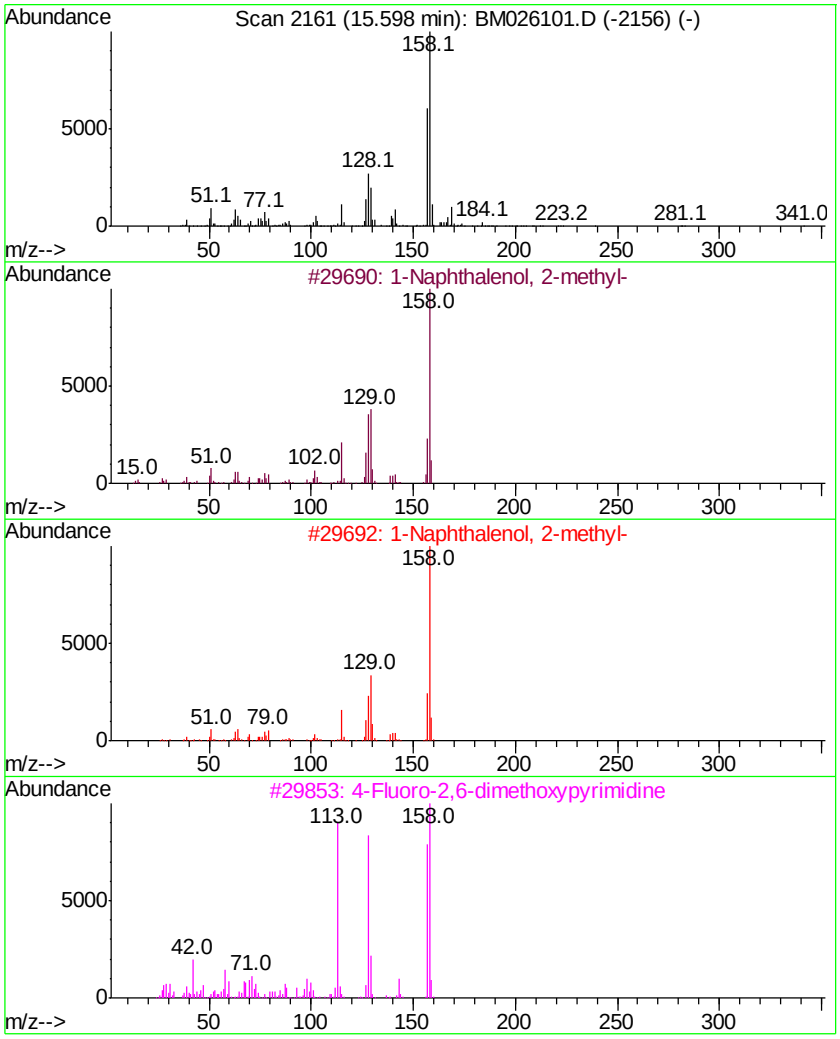
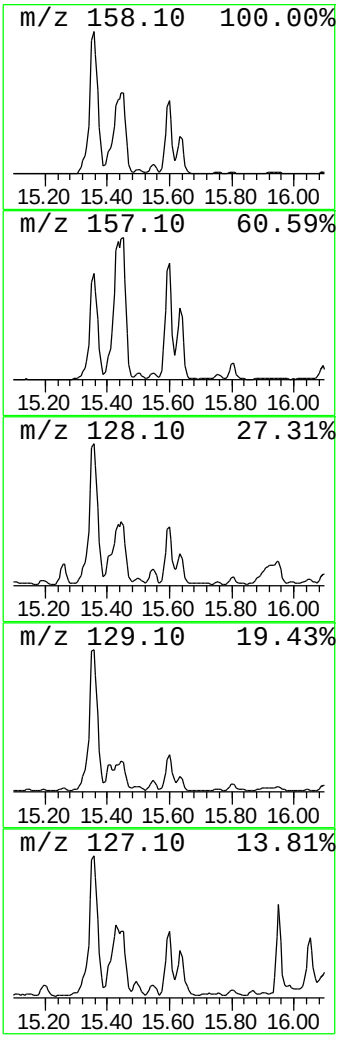
Peak Number 21 4-Fluoro-2,6-dimethoxypyrim... Concentration Rank 36

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.60	12.11 ng/ul	2186640	Phenanthrene-d10	16.89

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Naphthalenol, 2-methyl-	158	C11H10O	007469-77-4	62
2	1-Naphthalenol, 2-methyl-	158	C11H10O	007469-77-4	62
3	4-Fluoro-2,6-dimethoxypyrimidine	158	C6H7FN2O2	000658-87-7	59
4	7-Methyl-1-naphthol	158	C11H10O	006939-33-9	52
5	1H-Pyrazole, 3-methyl-1-phenyl-	158	C10H10N2	001128-54-7	50





Library Search Compound Report

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Operator : MA/SJ
Sample : L2737-13
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ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9B10

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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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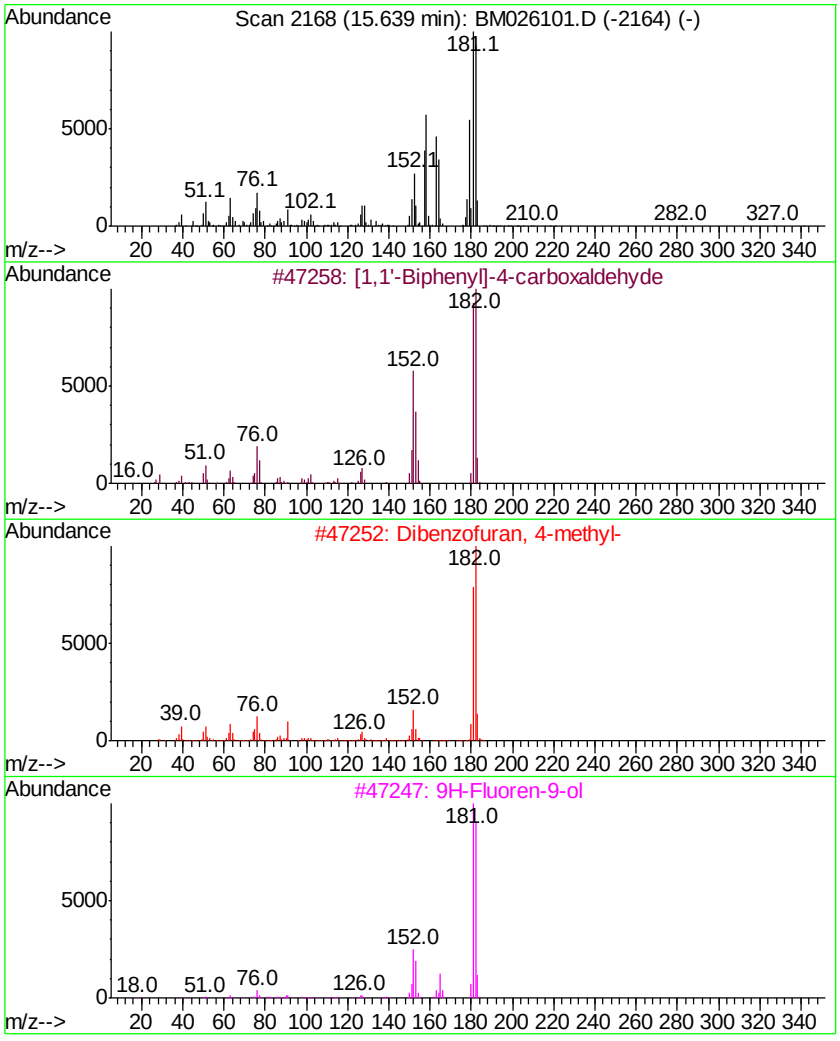
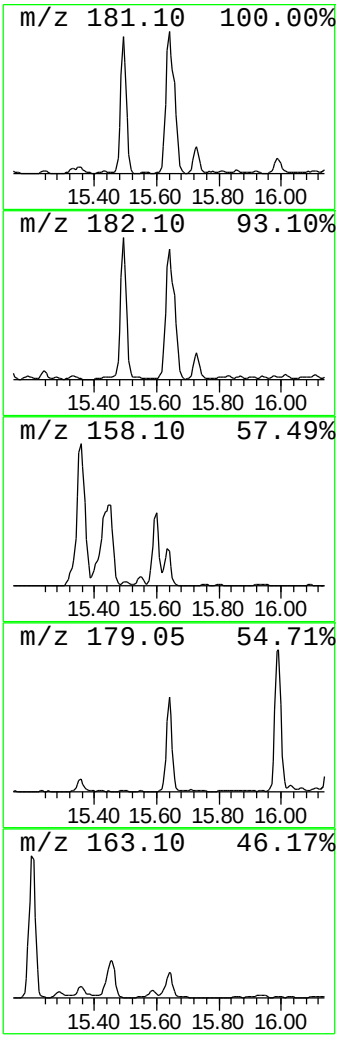
Peak Number 22 [1,1'-Biphenyl]-4-carboxald... Concentration Rank 33

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.64	16.23 ng/ul	2930810	Phenanthrene-d10	16.89

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	[1,1'-Biphenyl]-4-carboxaldehyde	182	C13H10O	003218-36-8	56
2	Dibenzofuran, 4-methyl-	182	C13H10O	007320-53-8	50
3	9H-Fluoren-9-ol	182	C13H10O	001689-64-1	43
4	9H-Fluoren-9-ol	182	C13H10O	001689-64-1	41
5	Norharmane, N-methyl-	182	C12H10N2	1000374-78-6	38





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ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9B10

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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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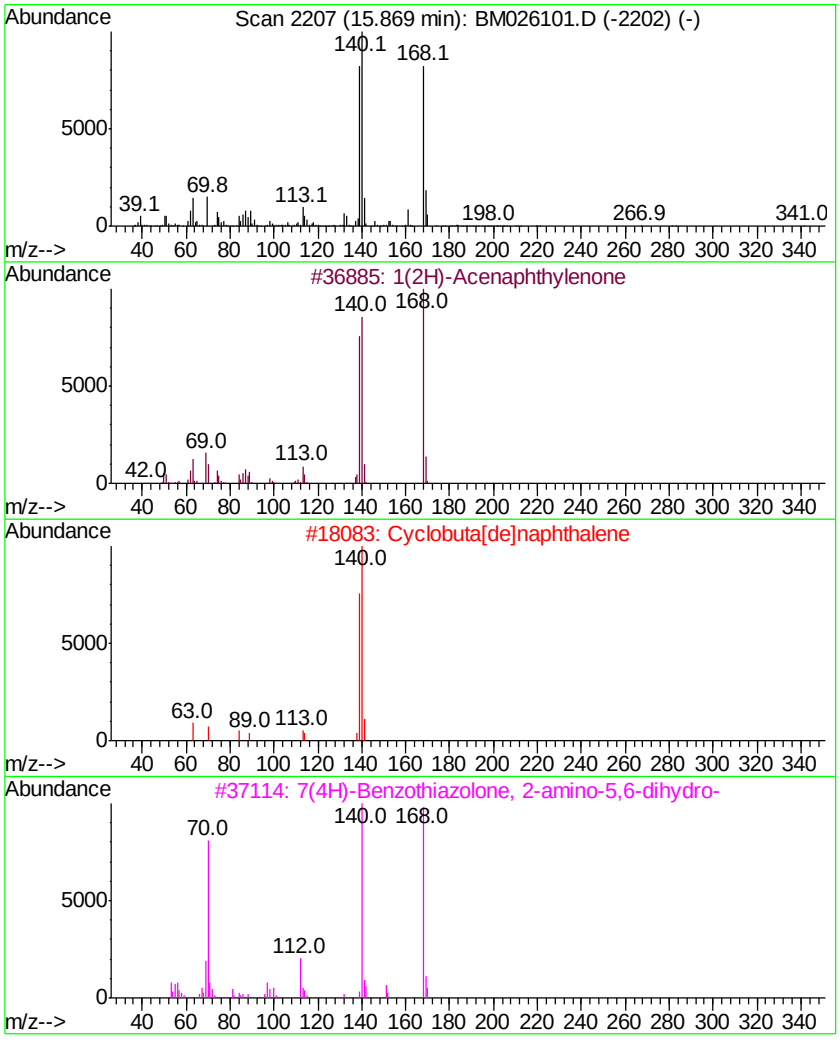
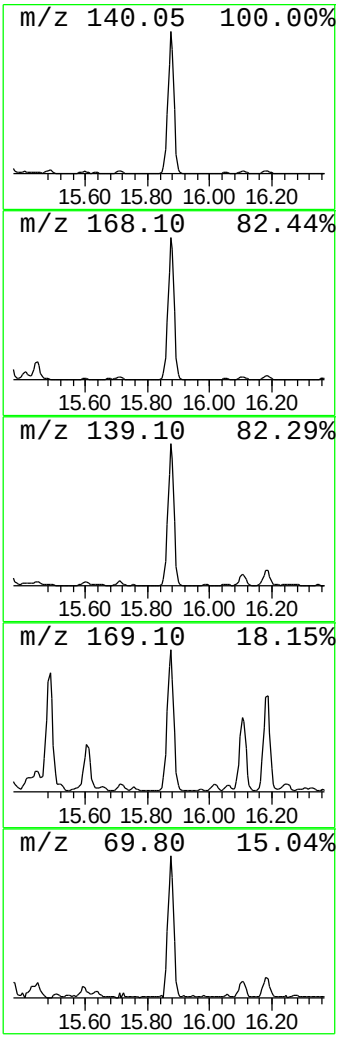
Peak Number 23 1(2H)-Acenaphthylenone Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.87	41.06 ng/ul	7416310	Phenanthrene-d10	16.89

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1(2H)-Acenaphthylenone	168	C12H8O	002235-15-6	96
2	Cyclobuta[de]naphthalene	140	C11H8	024973-91-9	52
3	7(4H)-Benzothiazolone, 2-amino-5...	168	C7H8N2OS	017583-10-7	50
4	Thiophene, 2,5-dipropyl-	168	C10H16S	054411-07-3	49
5	Oxidopamine	169	C8H11NO3	001199-18-4	47





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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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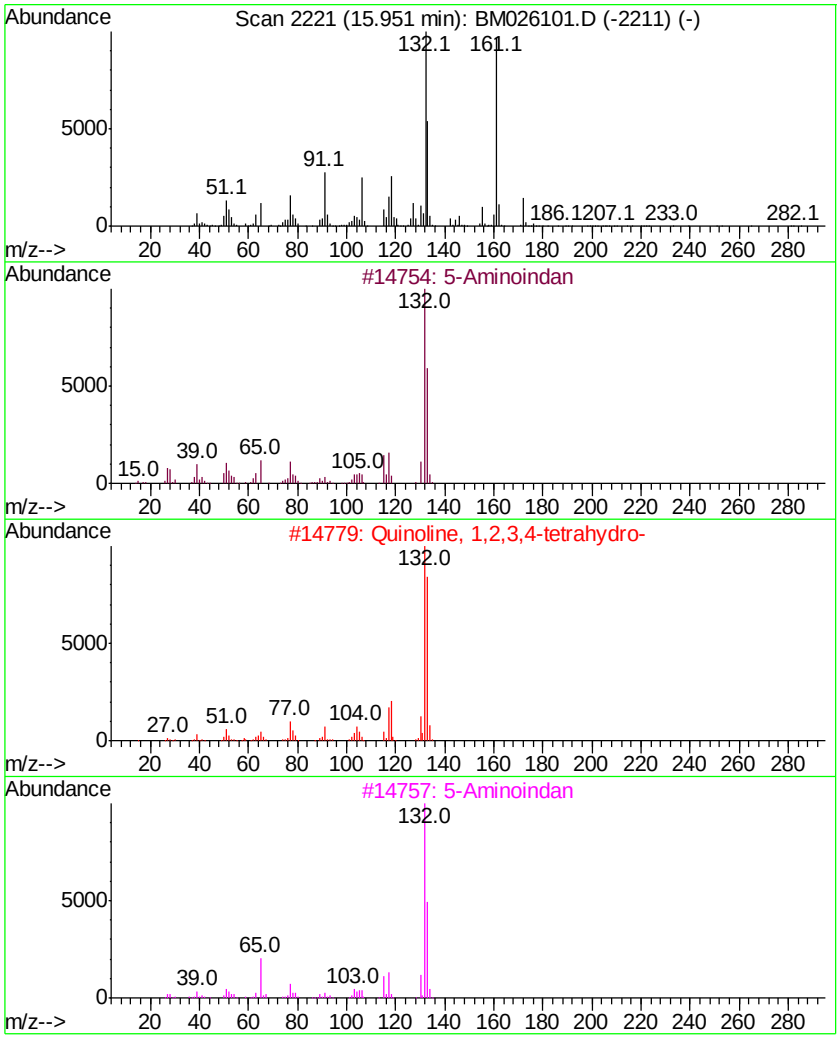
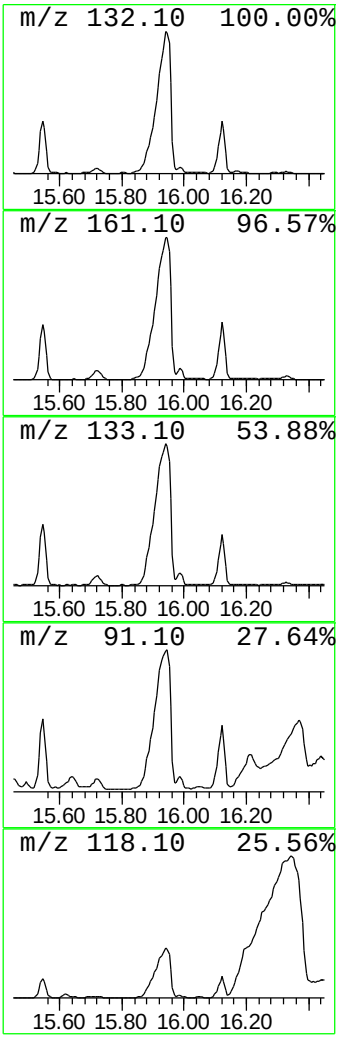
Peak Number 24 5-Aminoindan Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.95	82.79 ng/ul	14951400	Phenanthrene-d10	16.89

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5-Aminoindan	133	C9H11N	024425-40-9	86
2	Quinoline, 1,2,3,4-tetrahydro-	133	C9H11N	000635-46-1	46
3	5-Aminoindan	133	C9H11N	024425-40-9	45
4	Tranylcypromine	133	C9H11N	000155-09-9	43
5	5-Aminoindan	133	C9H11N	024425-40-9	43





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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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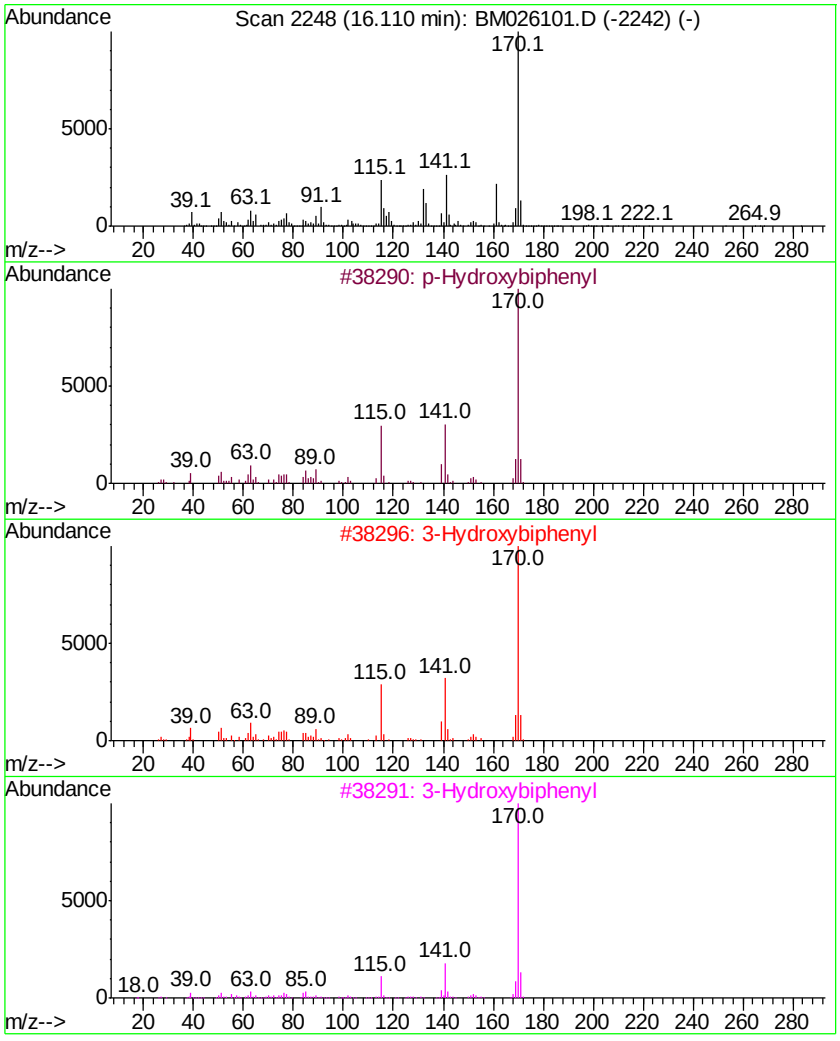
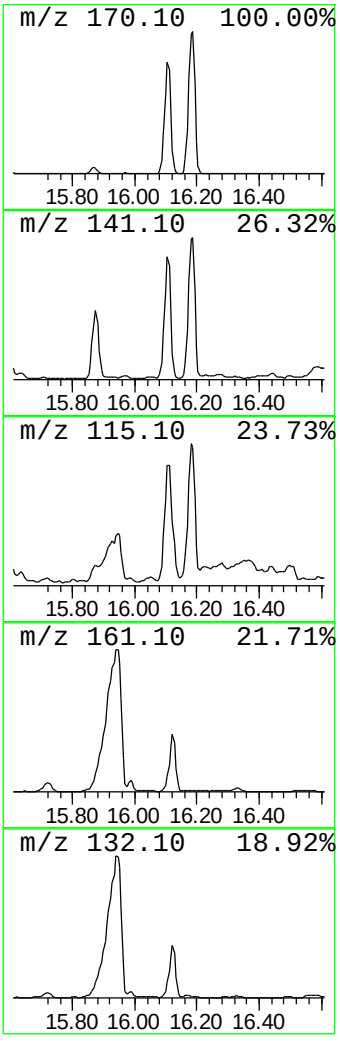
Peak Number 25 p-Hydroxybiphenyl Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.11	35.07 ng/ul	6333060	Phenanthrene-d10	16.89

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	p-Hydroxybiphenyl	170	C12H10O	000092-69-3	89
2	3-Hydroxybiphenyl	170	C12H10O	000580-51-8	89
3	3-Hydroxybiphenyl	170	C12H10O	000580-51-8	89
4	p-Hydroxybiphenyl	170	C12H10O	000092-69-3	86
5	3-Hydroxybiphenyl	170	C12H10O	000580-51-8	70





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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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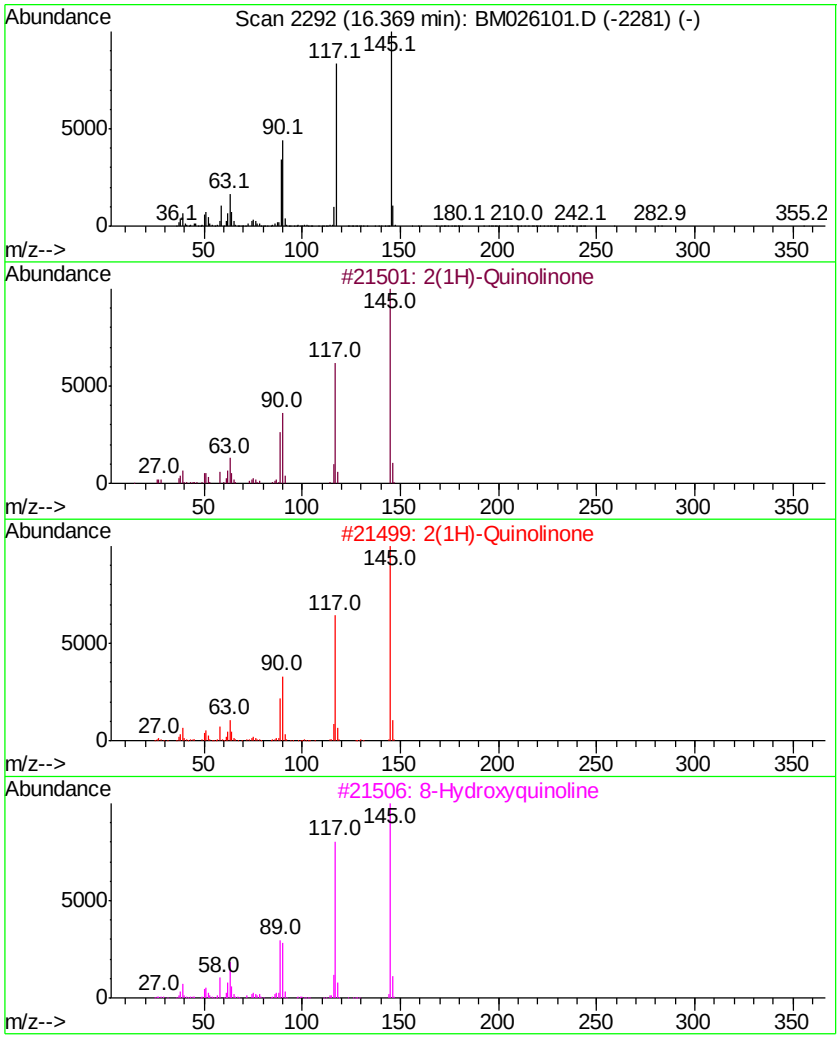
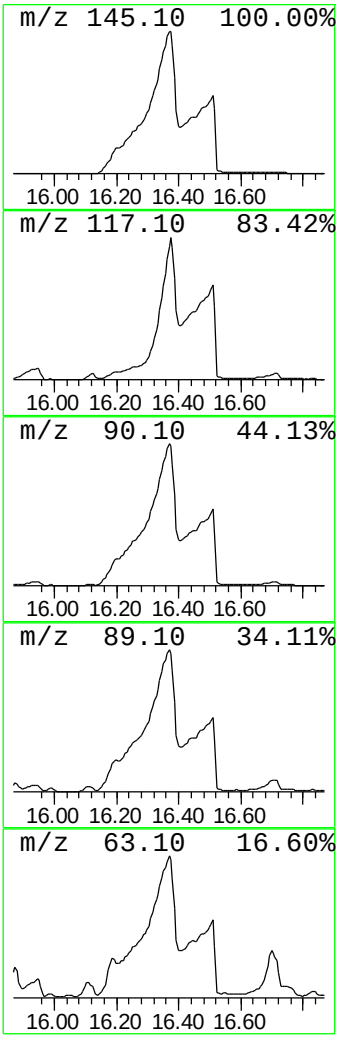
Peak Number 27 2(1H)-Quinolinone Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.37	147.12 ng/ul	26569600	Phenanthrene-d10	16.89

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2(1H)-Quinolinone	145	C9H7NO	000059-31-4	97
2	2(1H)-Quinolinone	145	C9H7NO	000059-31-4	95
3	8-Hydroxyquinoline	145	C9H7NO	000148-24-3	94
4	1(2H)-Isoquinolinone	145	C9H7NO	000491-30-5	93
5	8-Hydroxyquinoline	145	C9H7NO	000148-24-3	93





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Sample : L2737-13
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9B10

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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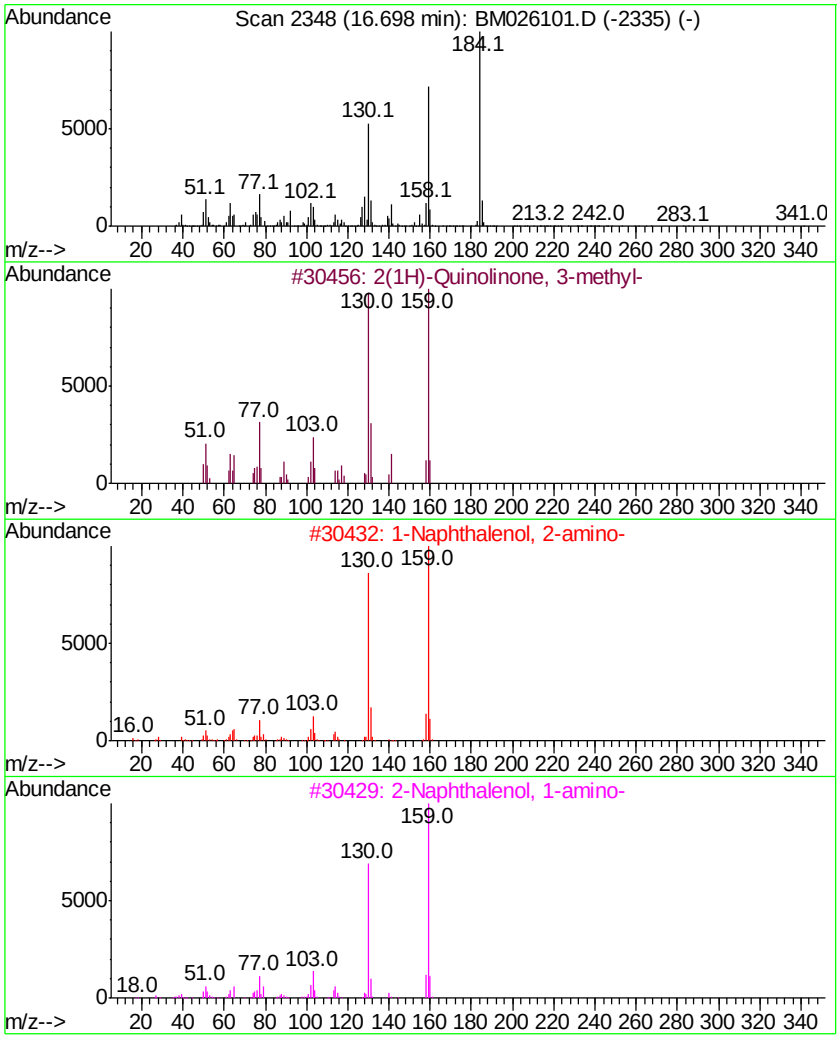
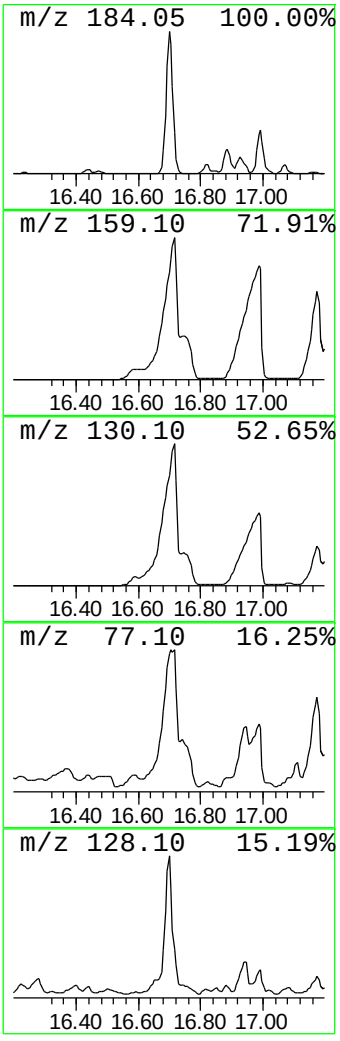
Peak Number 28 2(1H)-Quinolinone, 3-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.70	124.61 ng/ul	22504000	Phenanthrene-d10	16.89

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2(1H)-Quinolinone, 3-methyl-	159	C10H9NO	002721-59-7	91
2	1-Naphthalenol, 2-amino-	159	C10H9NO	000606-41-7	90
3	2-Naphthalenol, 1-amino-	159	C10H9NO	002834-92-6	78
4	2-Naphthalenol, 1-amino-	159	C10H9NO	002834-92-6	60
5	5-Amino-1-naphthol	159	C10H9NO	000083-55-6	55





Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM052220\
Data File : BM026101.D
Acq On : 23 May 2020 04:53
Operator : MA/SJ
Sample : L2737-13
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9B10

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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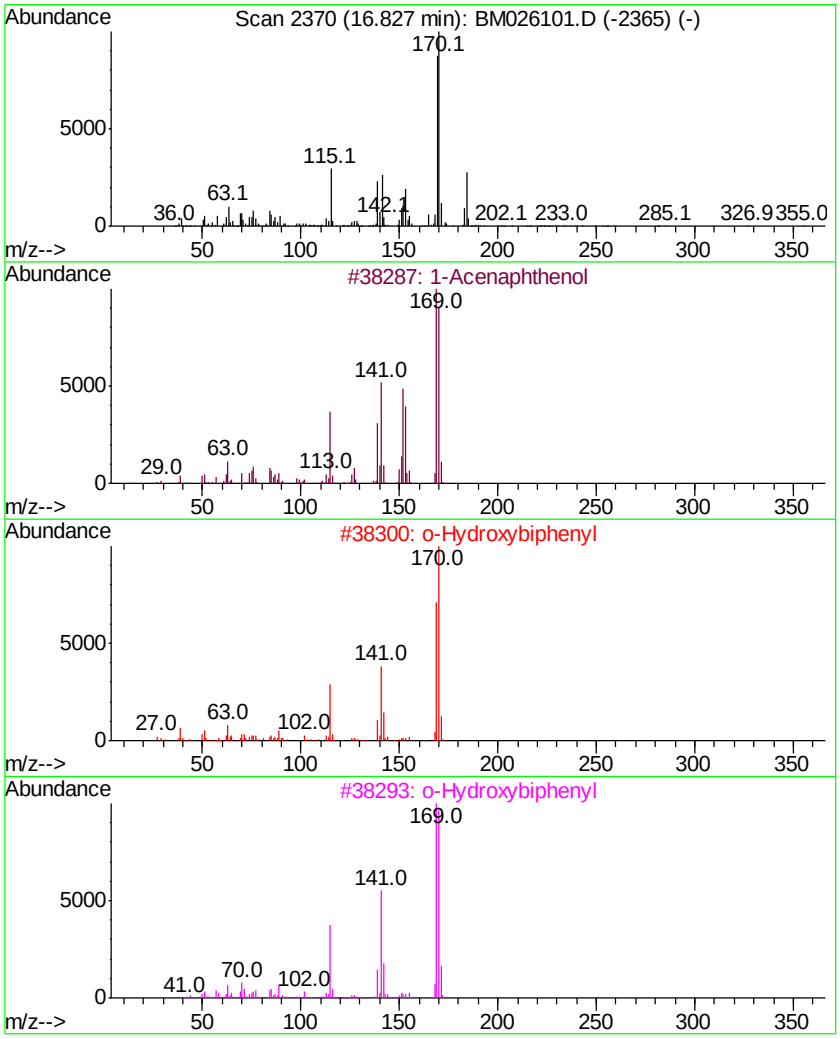
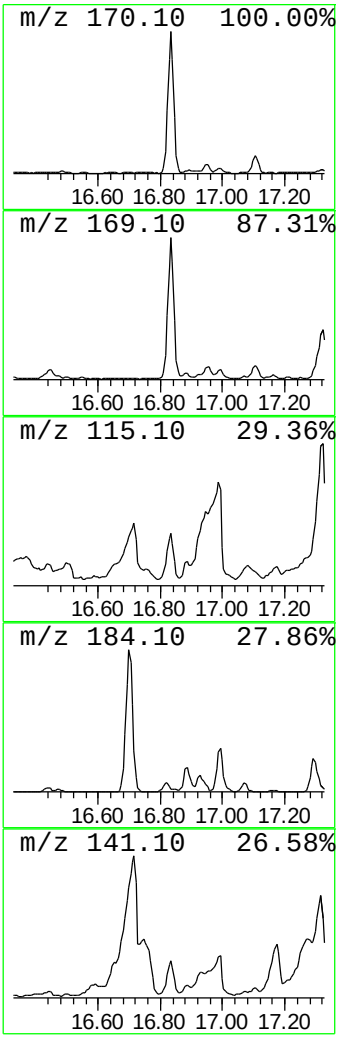
Peak Number 29 1-Acenaphthenol Concentration Rank 40

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.83	11.26 ng/ul	2034470	Phenanthrene-d10	16.89

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Acenaphthenol	170	C12H10O	006306-07-6	93
2	o-Hydroxybiphenyl	170	C12H10O	000090-43-7	74
3	o-Hydroxybiphenyl	170	C12H10O	000090-43-7	64
4	1-Acenaphthenol	170	C12H10O	006306-07-6	64
5	o-Hydroxybiphenyl	170	C12H10O	000090-43-7	58





Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM052220\
Data File : BM026101.D
Acq On : 23 May 2020 04:53
Operator : MA/SJ
Sample : L2737-13
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9B10

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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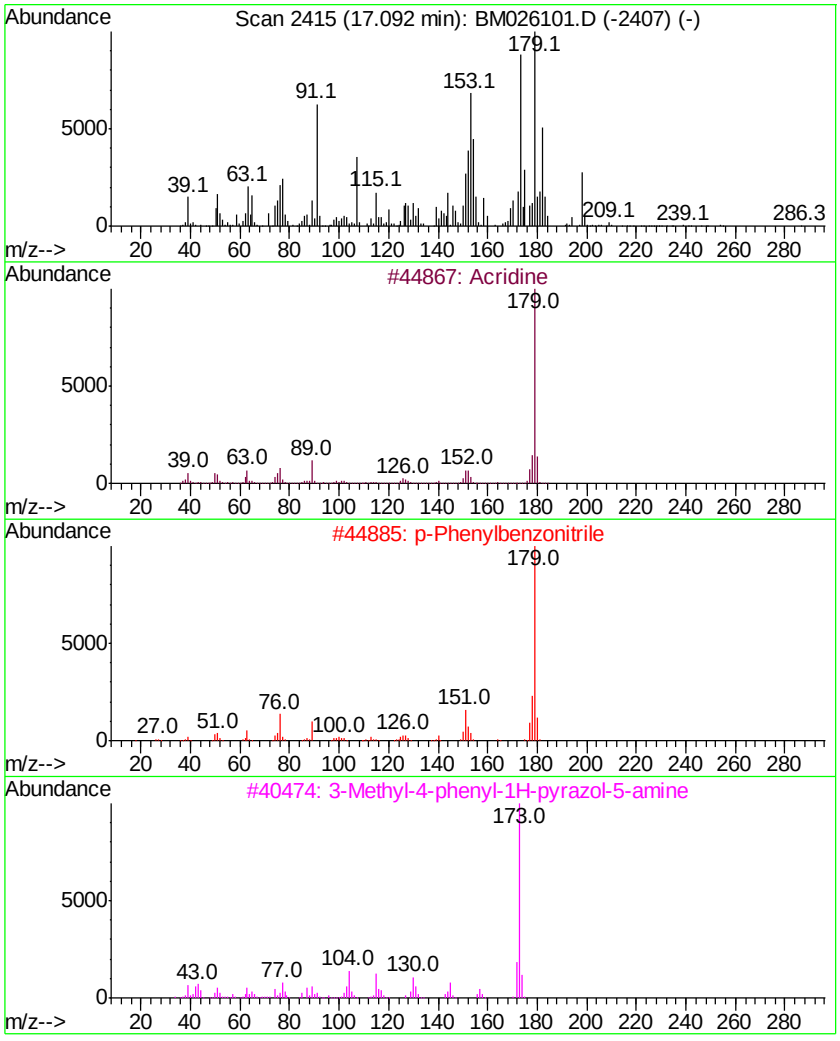
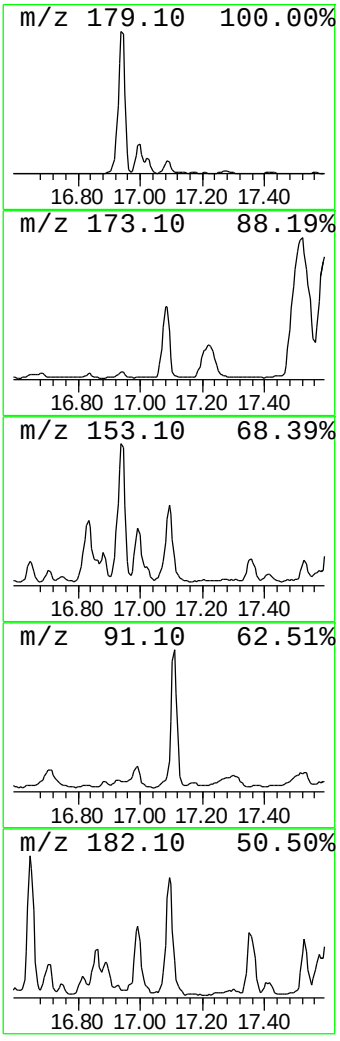
Peak Number 30 unknown-03 Concentration Rank 43

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.09	10.01 ng/ul	1808030	Phenanthrene-d10	16.89

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Acridine	179	C13H9N	000260-94-6	38
2	p-Phenylbenzonitrile	179	C13H9N	002920-38-9	35
3	3-Methyl-4-phenyl-1H-pyrazol-5-a...	173	C10H11N3	060419-81-0	35
4	Benzo[g]quinoline	179	C13H9N	000260-36-6	30
5	Acridine	179	C13H9N	000260-94-6	30





Tentatively Identified Compound (LSC) summary

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM052220\
Data File : BM026101.D
Acq On : 23 May 2020 04:53
Operator : MA/SJ
Sample : L2737-13
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9B10

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM051320MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

					Internal Standard			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Benzofuran	7.22	15.8	ng/ul	2724530	1	7.49	3449800	20.0
Eucalyptol	7.80	25.7	ng/ul	4435280	1	7.49	3449800	20.0
Benzene, 1-propenyl-	7.86	67.8	ng/ul	11693100	1	7.49	3449800	20.0
Indene	8.02	21.0	ng/ul	3625080	1	7.49	3449800	20.0
Bicyclo[2.2.1]hep...	8.73	22.5	ng/ul	3887560	1	7.49	3449800	20.0
Cyclohexanemethan...	9.69	2.1	ng/ul	21267200	2	10.29	203457000	20.0
Phenol, 3,5-dimet...	9.83	2.7	ng/ul	27857700	2	10.29	203457000	20.0
1H-Inden-5-ol, 2,...	12.34	21.0	ng/ul	3805800	3	14.14	3625730	20.0
unknown-01	12.87	8.5	ng/ul	1534060	3	14.14	3625730	20.0
Naphthalene, 2,7-...	13.29	15.1	ng/ul	2729630	3	14.14	3625730	20.0
Naphthalene, 2,3-...	13.46	23.7	ng/ul	4293960	3	14.14	3625730	20.0
Naphthalene, 1,7-...	13.51	12.1	ng/ul	2185560	3	14.14	3625730	20.0
Naphthalene, 1,3-...	13.69	8.8	ng/ul	1593840	3	14.14	3625730	20.0
2-Naphthalenecarb...	14.27	19.9	ng/ul	3603210	3	14.14	3625730	20.0
o-Hydroxybiphenyl	14.43	46.6	ng/ul	8455480	3	14.14	3625730	20.0
1,3-Phenylene, bi...	14.65	20.1	ng/ul	3648340	3	14.14	3625730	20.0
2-Naphthalenamine...	15.02	281.2	ng/ul	50980100	3	14.14	3625730	20.0
1-Naphthalenol, 2...	15.36	32.9	ng/ul	5964880	3	14.14	3625730	20.0
unknown-02	15.45	32.2	ng/ul	5830460	3	14.14	3625730	20.0
1(2H)-Quinolineca...	15.55	18.1	ng/ul	3274700	4	16.89	3612060	20.0
4-Fluoro-2,6-dime...	15.60	12.1	ng/ul	2186640	4	16.89	3612060	20.0
[1,1'-Biphenyl]-4...	15.64	16.2	ng/ul	2930810	4	16.89	3612060	20.0
1(2H)-Acenaphthyl...	15.87	41.1	ng/ul	7416310	4	16.89	3612060	20.0
5-Aminoindan	15.95	82.8	ng/ul	14951400	4	16.89	3612060	20.0
p-Hydroxybiphenyl	16.11	35.1	ng/ul	6333060	4	16.89	3612060	20.0
2(1H)-Quinolinone	16.37	147.1	ng/ul	26569600	4	16.89	3612060	20.0
2(1H)-Quinolinone...	16.70	124.6	ng/ul	22504000	4	16.89	3612060	20.0
1-Acenaphthenol	16.83	11.3	ng/ul	2034470	4	16.89	3612060	20.0
unknown-03	17.09	10.0	ng/ul	1808030	4	16.89	3612060	20.0