LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM052620\
Data File : BM026124.D
Acq On : 26 May 2020 21:51
Operator : MA/SJ
Sample : L2737-12DL 10X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9B27DL

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	7.210	729	735	744	rVB	399676	621015	0.66%	0.303%
2	7.487	773	782	789	rBV	687239	1135543	1.21%	0.553%
3	7.798	828	835	839	rBV	151623	236472	0.25%	0.115%
4	7.857	839	845	853	rVV	2531171	3791354	4.05%	1.847%
5	7.939	853	859	865	rVV	298642	446785	0.48%	0.218%
6	8.016	865	872	880	rVB	1138309	1763107	1.89%	0.859%
7	8.263	909	914	919	rVV	202357	329568	0.35%	0.161%
8	8.322	919	924	933	rVB	387179	636743	0.68%	0.310%
9	8.634	972	977	979	rBV	82308	136356	0.15%	0.066%
10	8.716	986	991	998	rVB	261828	395233	0.42%	0.193%
11	8.828	1003	1010	1016	rBV	83165	134680	0.14%	0.066%
12	8.904	1016	1023	1025	rBV	105580	167492	0.18%	0.082%
13	8.951	1025	1031	1037	rVV	217720	477445	0.51%	0.233%
14	9.457	1111	1117	1120	rBV	486289	742813	0.79%	0.362%
15	9.492	1120	1123	1130	rVV3	258805	465617	0.50%	0.227%
16	9.639	1141	1148	1151	rBV	1118357	1790104	1.91%	0.872%
17	9.675	1151	1154	1165	rVB3	589130	974834	1.04%	0.475%
18	9.769	1165	1170	1184	rVB	822374	1423336	1.52%	0.693%
19	9.892	1184	1191	1195	rBV	223971	420707	0.45%	0.205%
20	10.151	1230	1235	1240	rBV	148216	213869	0.23%	0.104%
21	10.339	1254	1267	1272	rVB2	34997983	93517627	100.00%	45.555%
22	10.422	1274	1281	1292	rVB	2678805	4377874	4.68%	2.133%
23	10.622	1308	1315	1321	rBV4	68385	166202	0.18%	0.081%
24	10.804	1342	1346	1350	rBV	105518	158587	0.17%	0.077%
25	11.098	1389	1396	1403	rBV2	293836	477236	0.51%	0.232%
26	11.286	1420	1428	1433	rBV	90655	150912	0.16%	0.074%
27	11.345	1433	1438	1452	rVB2	143469	354075	0.38%	0.172%
28	11.639	1481	1488	1494	rBV	211453	390475	0.42%	0.190%
29	11.922	1526	1536	1543	rBV	3771740	5917091	6.33%	2.882%
30	12.033	1550	1555	1561	rVB3	173193	300823	0.32%	0.147%
31	12.145	1561	1574	1584	rVB	2176040	3394439	3.63%	1.654%
32	12.327	1600	1605	1610	rVV2	125510	205326	0.22%	0.100%
33	12.957	1701	1712	1723	rVB	698383	1346093	1.44%	0.656%
34	13.157	1741	1746	1759	rVB	148753	331049	0.35%	0.161%

LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM052620\
Data File : BM026124.D
Acq On : 26 May 2020 21:51
Operator : MA/SJ
Sample : L2737-12DL 10X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9B27DL

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	13.286	1759	1768	1770	rBV3	153493	297130	0.32%	0.145%
36	13.451	1791	1796	1801	rVV	256903	465396	0.50%	0.227%
37	13.504	1801	1805	1809	rVV	158387	239604	0.26%	0.117%
38	13.557	1809	1814	1818	rVB	179587	261967	0.28%	0.128%
39	13.686	1832	1836	1840	rVV	113470	173645	0.19%	0.085%
40	13.821	1851	1859	1861	rBV	220032	347056	0.37%	0.169%
41	13.851	1861	1864	1872	rVB2	182216	278980	0.30%	0.136%
42	14.133	1904	1912	1917	rBV	1247454	1905069	2.04%	0.928%
43	14.198	1917	1923	1929	rVV	3399819	4760963	5.09%	2.319%
44	14.263	1929	1934	1939	rVV	802630	1108619	1.19%	0.540%
45	14.327	1939	1945	1954	rVV	582371	885692	0.95%	0.431%
46	14.416	1954	1960	1970	rVB2	797783	1519327	1.62%	0.740%
47	14.533	1970	1980	1989	rVB	1814741	2949986	3.15%	1.437%
48	14.633	1991	1997	2008	rVB2	115677	230566	0.25%	0.112%
49	14.851	2029	2034	2035	rVV	224658	264010	0.28%	0.129%
50	14.874	2035	2038	2043	rVV	406803	550237	0.59%	0.268%
51	15.010	2056	2061	2067	rBV	134705	224022	0.24%	0.109%
52	15.133	2077	2082	2086	rVB	296123	412446	0.44%	0.201%
53	15.186	2086	2091	2099	rVB	1684391	2240512	2.40%	1.091%
54	15.268	2099	2105	2109	rBV4	79005	157727	0.17%	0.077%
55	15.333	2109	2116	2123	rVV	487127	1123874	1.20%	0.547%
56	15.427	2123	2132	2138	rVV4	340144	896119	0.96%	0.437%
57	15.504	2138	2145	2153	rVV3	335350	629246	0.67%	0.307%
58	15.574	2153	2157	2161	rVV	186726	289298	0.31%	0.141%
59	15.621	2161	2165	2175	rVB4	173404	431643	0.46%	0.210%
60	15.880	2200	2209	2216	rBV	638470	1426437	1.53%	0.695%
61	16.157	2246	2256	2267	rVB4	2510294	9838844	10.52%	4.793%
62	16.251	2268	2272	2285	rVB2	103448	263869	0.28%	0.129%
63	16.504	2303	2315	2319	rVV	1286474	2897211	3.10%	1.411%
64	16.580	2319	2328	2333	rVB	874730	1942241	2.08%	0.946%
65	16.668	2339	2343	2352	rVB	688114	1073071	1.15%	0.523%
66	16.768	2352	2360	2363	rBV	693596	1311476	1.40%	0.639%
67	16.804	2363	2366	2371	rVB2	187443	261407	0.28%	0.127%
68	16.880	2371	2379	2383	rBV	1479957	2354817	2.52%	1.147%
69	16.921	2383	2386	2390	rVB	1874929	2261217	2.42%	1.102%
70	16.980	2390	2396	2399	rBV3	681379	1210265	1.29%	0.590%
71	17.021	2399	2403	2406	rBV3	725646	1257824	1.35%	0.613%

LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM052620\
Data File : BM026124.D
Acq On : 26 May 2020 21:51
Operator : MA/SJ
Sample : L2737-12DL 10X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9B27DL

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

72	17.068	2407	2411	2413	rBV	465094	661769	0.71%	0.322%
73	17.127	2419	2421	2427	rVB	364478	514435	0.55%	0.251%
74	17.198	2427	2433	2438	rVB	447912	616912	0.66%	0.301%
75	17.292	2438	2449	2456	rBV	4795286	7548599	8.07%	3.677%
76	17.392	2456	2466	2471	rBV2	1226551	2119515	2.27%	1.032%
77	17.456	2471	2477	2485	rVB	1305015	2129127	2.28%	1.037%
78	17.551	2491	2493	2500	rVB2	95577	177790	0.19%	0.087%
79	17.639	2505	2508	2512	rVB2	196909	233869	0.25%	0.114%
80	17.727	2517	2523	2527	rBV2	441852	736018	0.79%	0.359%
81	17.804	2532	2536	2539	rVV	524665	714898	0.76%	0.348%
82	17.833	2539	2541	2545	rVV	209390	247375	0.26%	0.121%
83	17.886	2545	2550	2555	rVB	760785	1015226	1.09%	0.495%
84	17.951	2556	2561	2568	rBV3	144775	277230	0.30%	0.135%
85	18.051	2573	2578	2583	rBV3	231208	452907	0.48%	0.221%
86	18.104	2585	2587	2593	rVV2	116483	199157	0.21%	0.097%
87	18.215	2600	2606	2611	rVB3	130732	297759	0.32%	0.145%
88	18.268	2611	2615	2620	rBV	182551	253435	0.27%	0.123%
89	18.945	2725	2730	2738	rVV	442559	607874	0.65%	0.296%
90	19.145	2760	2764	2771	rBV2	156466	331617	0.35%	0.162%
91	19.280	2782	2787	2790	rBV	481172	684655	0.73%	0.334%
92	19.309	2790	2792	2794	rVV	261677	311432	0.33%	0.152%
93	19.345	2794	2798	2804	rVB2	956325	1262161	1.35%	0.615%
94	19.698	2853	2858	2863	rBV2	267578	428605	0.46%	0.209%
95	19.768	2865	2870	2876	rVB	1082524	1508487	1.61%	0.735%
96	20.374	2969	2973	2977	rBV2	103785	174559	0.19%	0.085%
97	21.080	3087	3093	3097	rBV	2022895	2560844	2.74%	1.247%
98	21.544	3168	3172	3176	rBV	151313	205634	0.22%	0.100%
99	23.062	3426	3430	3436	rVB	278349	431984	0.46%	0.210%
100	23.197	3447	3453	3461	rVB	1436519	2448066	2.62%	1.193%

Sum of corrected areas: 205284634

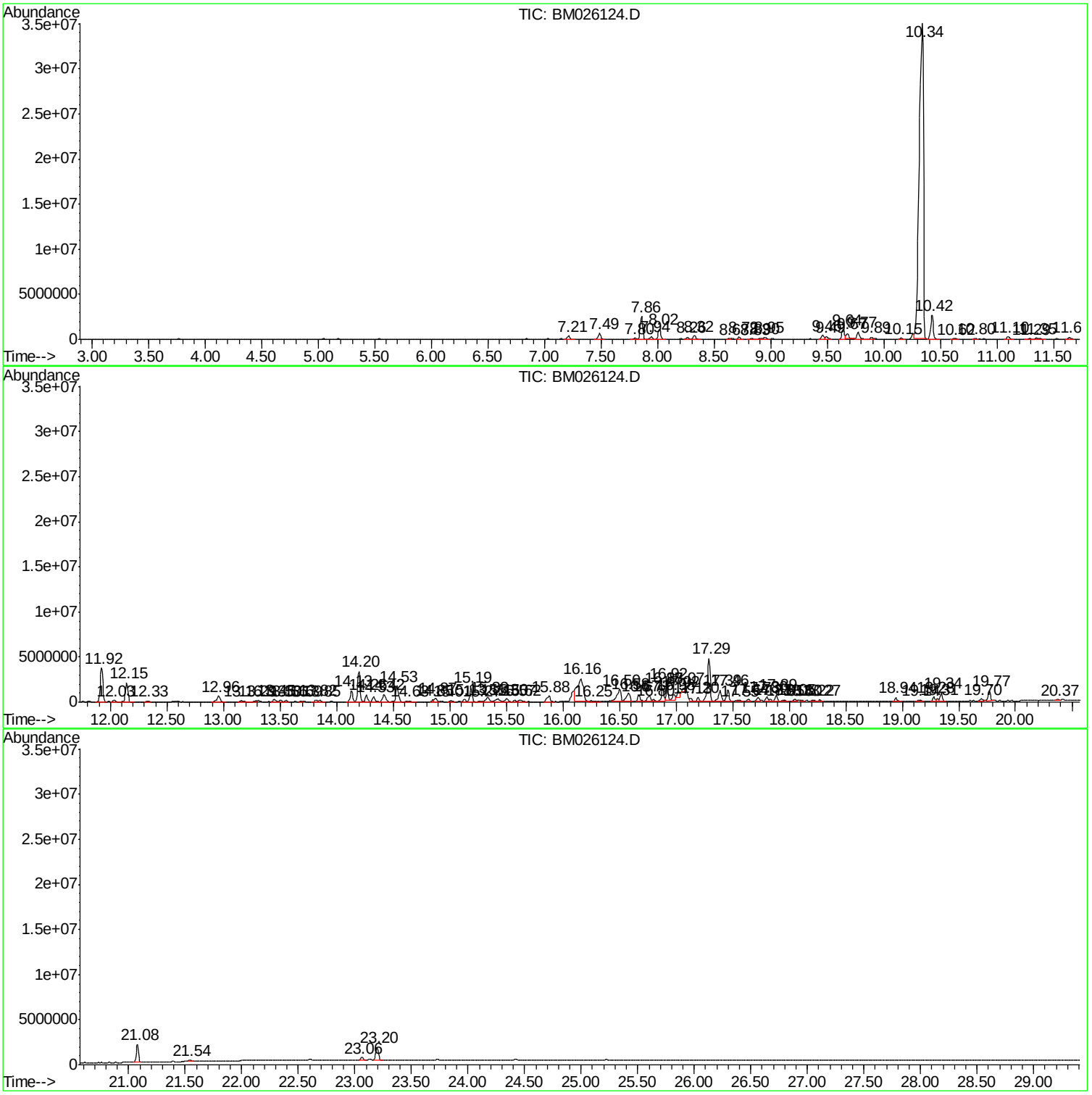
LSC Report - Integrated Chromatogram

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM052620\
Data File : BM026124.D
Acq On : 26 May 2020 21:51
Operator : MA/SJ
Sample : L2737-12DL 10X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
F9B27DL

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



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM052620\
Data File : BM026124.D
Acq On : 26 May 2020 21:51
Operator : MA/SJ
Sample : L2737-12DL 10X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9B27DL

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

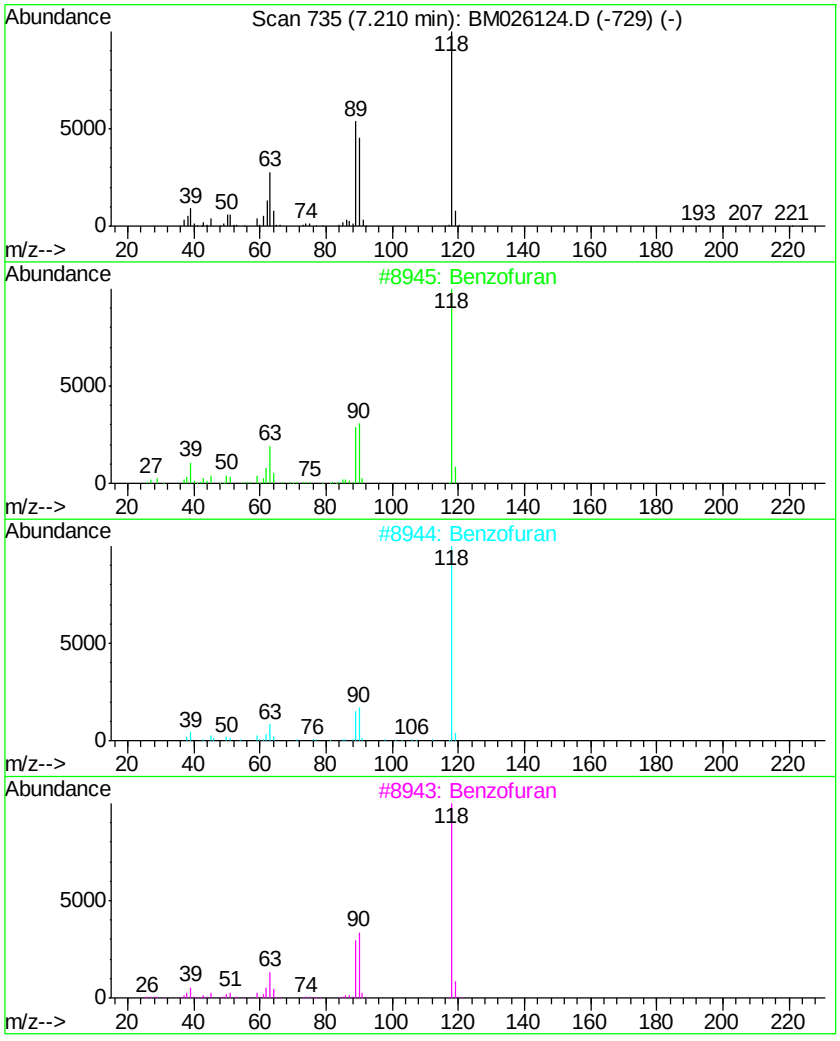
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

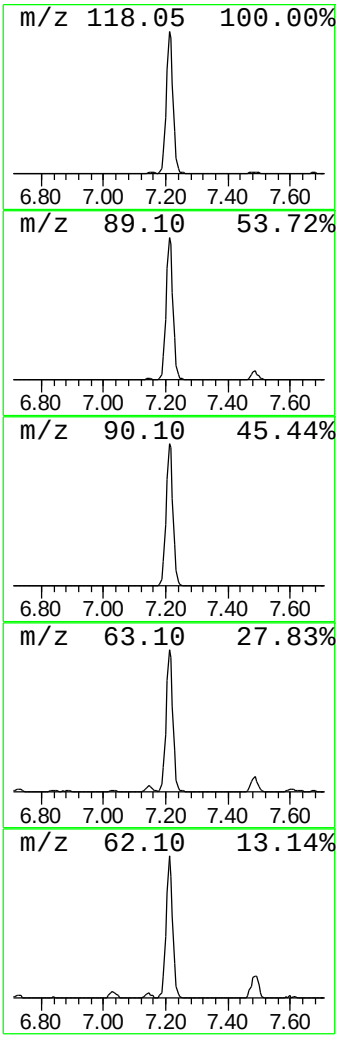
****

Peak Number 1 Benzofuran Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.21	10.94 ng/ul	621015	1,4-Dichlorobenzene-d4	7.49

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzofuran	118	C8H6O	000271-89-6	94
2	Benzofuran	118	C8H6O	000271-89-6	90
3	Benzofuran	118	C8H6O	000271-89-6	90
4	3-Hydroxyphenylacetylene	118	C8H6O	010401-11-3	72
5	3-Pyridineacetonitrile	118	C7H6N2	006443-85-2	38





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM052620\
Data File : BM026124.D
Acq On : 26 May 2020 21:51
Operator : MA/SJ
Sample : L2737-12DL 10X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9B27DL

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

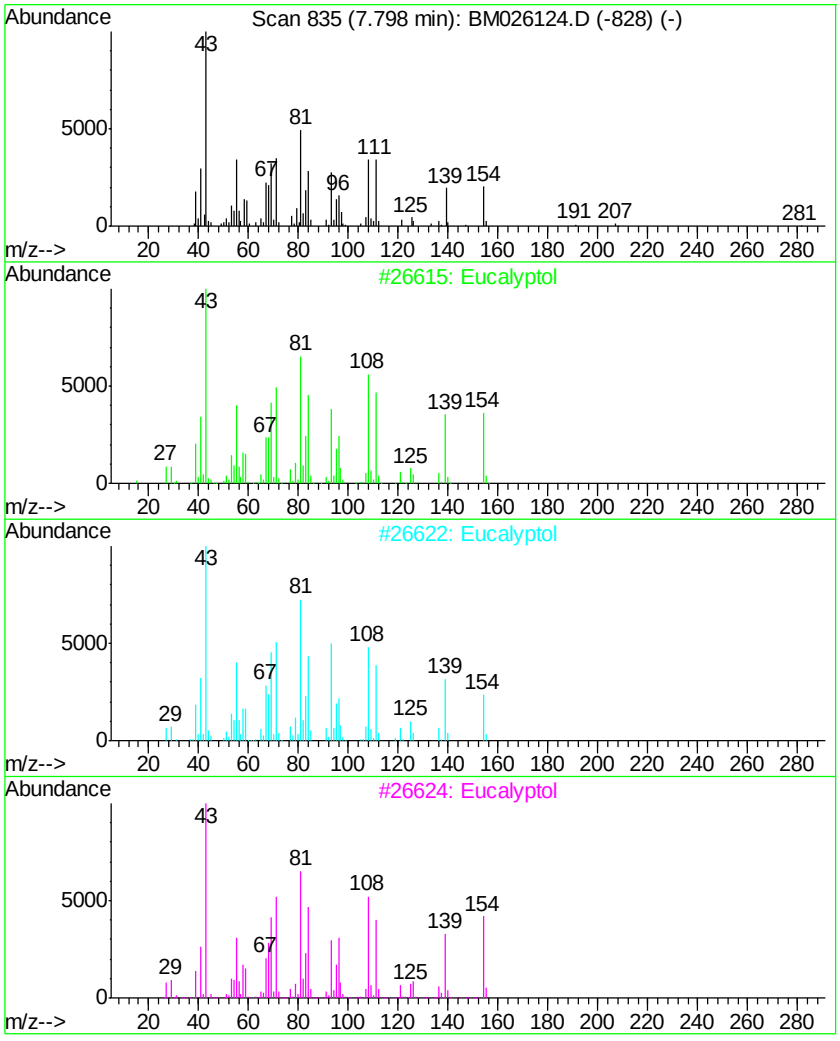
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

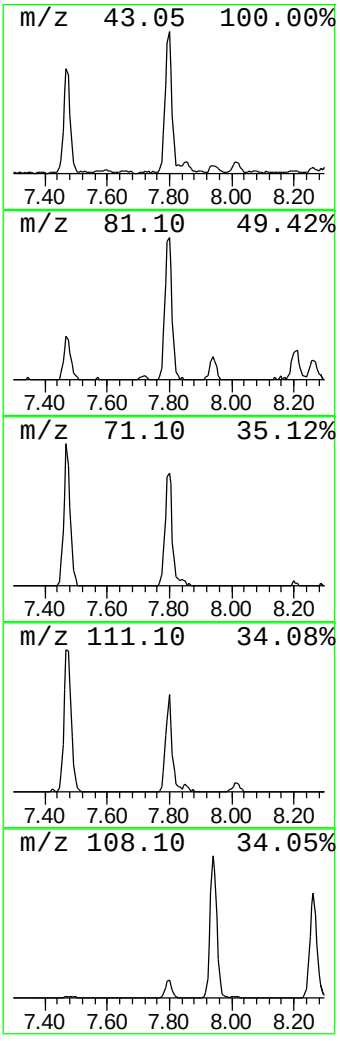
****

Peak Number 2 Eucalyptol Concentration Rank 29

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.80	4.16 ng/ul	236472	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	91
4	Eucalyptol	154	C10H18O	000470-82-6	91
5	Eucalyptol	154	C10H18O	000470-82-6	74





Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM052620\
Data File : BM026124.D
Acq On : 26 May 2020 21:51
Operator : MA/SJ
Sample : L2737-12DL 10X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9B27DL

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

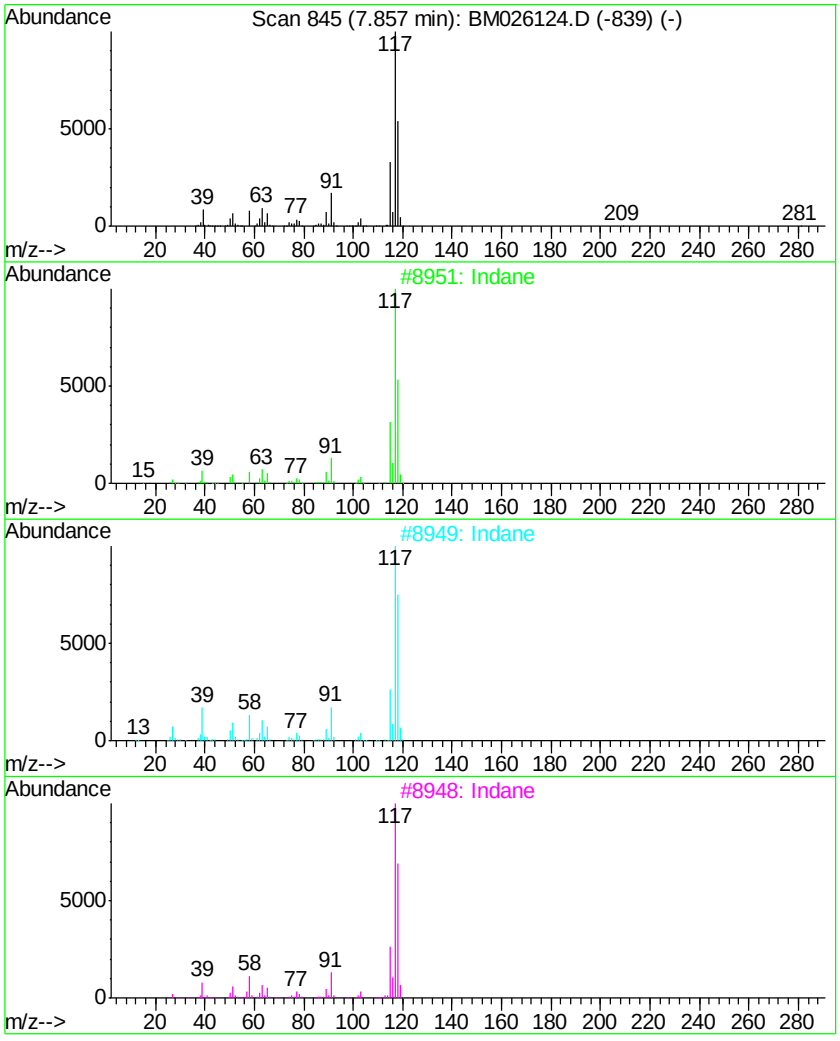
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

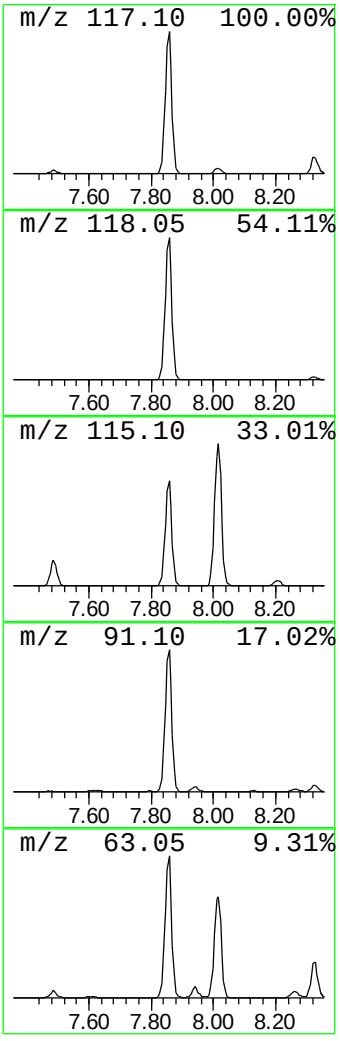
****

Peak Number 3 Tetracyclo[3.3.1.0(2,8).0(4... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.86	66.78 ng/ul	3791350	1,4-Dichlorobenzene-d4	7.49

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	93
2	Indane	118	C9H10	000496-11-7	93
3	Indane	118	C9H10	000496-11-7	81
4	Tetracyclo[3.3.1.0(2,8).0(4,6)]-...	118	C9H10	1000191-13-7	80
5	Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	60





Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM052620\
Data File : BM026124.D
Acq On : 26 May 2020 21:51
Operator : MA/SJ
Sample : L2737-12DL 10X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9B27DL

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

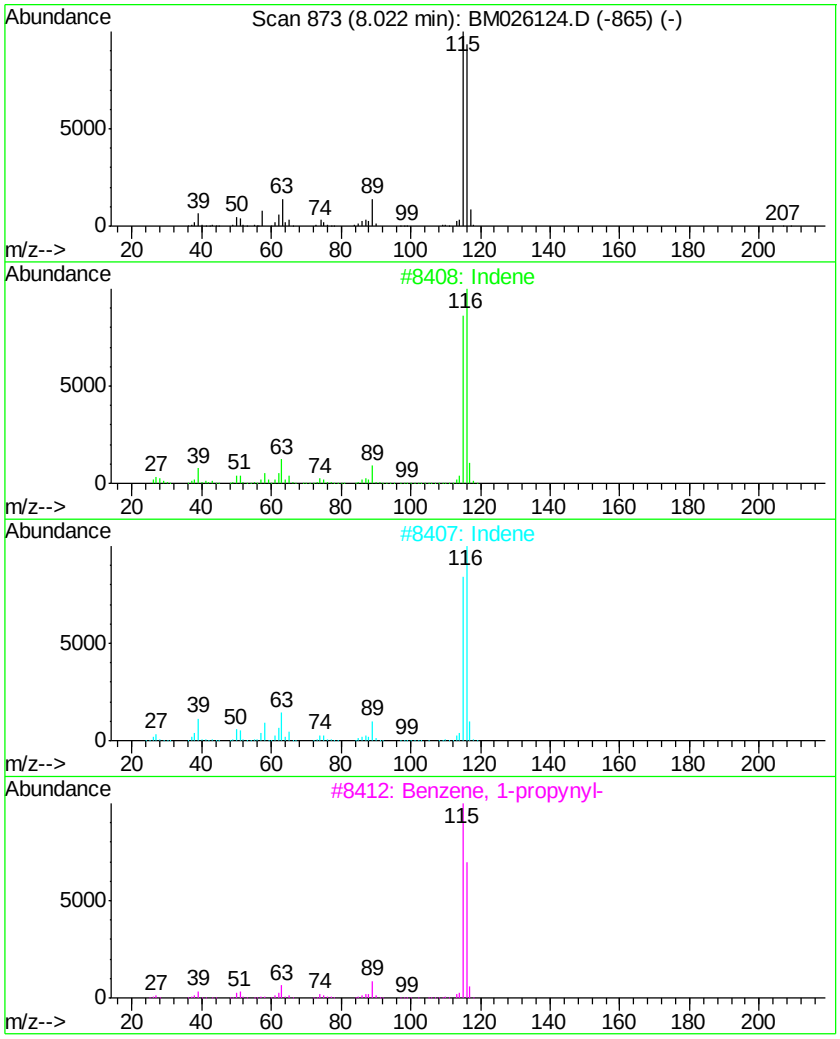
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

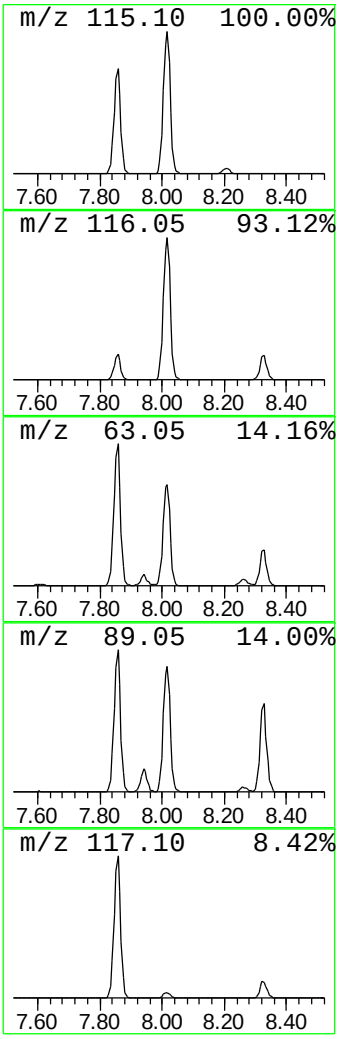
****

Peak Number 4 Indene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.02	31.05 ng/ul	1763110	1,4-Dichlorobenzene-d4	7.49

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indene	116	C9H8	000095-13-6	95
2	Indene	116	C9H8	000095-13-6	95
3	Benzene, 1-propynyl-	116	C9H8	000673-32-5	91
4	3-Methylphenylacetylene	116	C9H8	000766-82-5	91
5	Benzene, 1-propynyl-	116	C9H8	000673-32-5	91





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM052620\
Data File : BM026124.D
Acq On : 26 May 2020 21:51
Operator : MA/SJ
Sample : L2737-12DL 10X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9B27DL

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

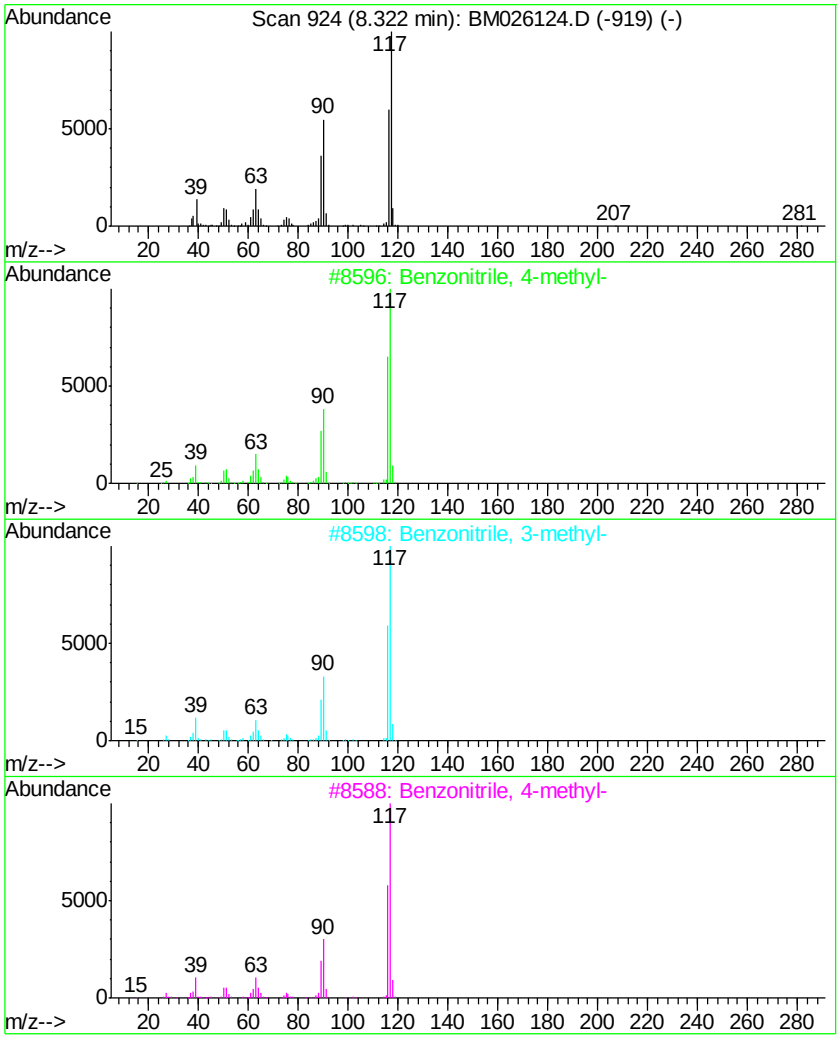
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

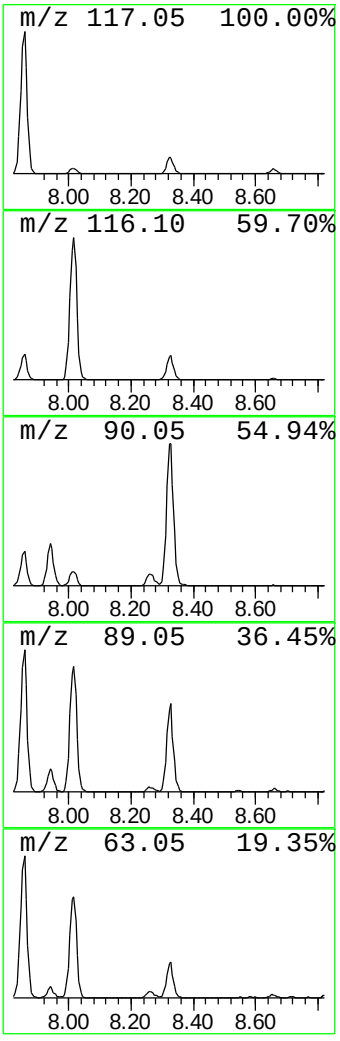
****

Peak Number 5 Benzonitrile, 4-methyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.32	11.21 ng/ul	636743	1,4-Dichlorobenzene-d4	7.49

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzonitrile, 4-methyl-	117	C8H7N	000104-85-8	97
2	Benzonitrile, 3-methyl-	117	C8H7N	000620-22-4	97
3	Benzonitrile, 4-methyl-	117	C8H7N	000104-85-8	97
4	Benzonitrile, 2-methyl-	117	C8H7N	000529-19-1	96
5	Benzonitrile, 4-methyl-	117	C8H7N	000104-85-8	96





Library Search Compound Report

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Data File : BM026124.D
Acq On : 26 May 2020 21:51
Operator : MA/SJ
Sample : L2737-12DL 10X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9B27DL

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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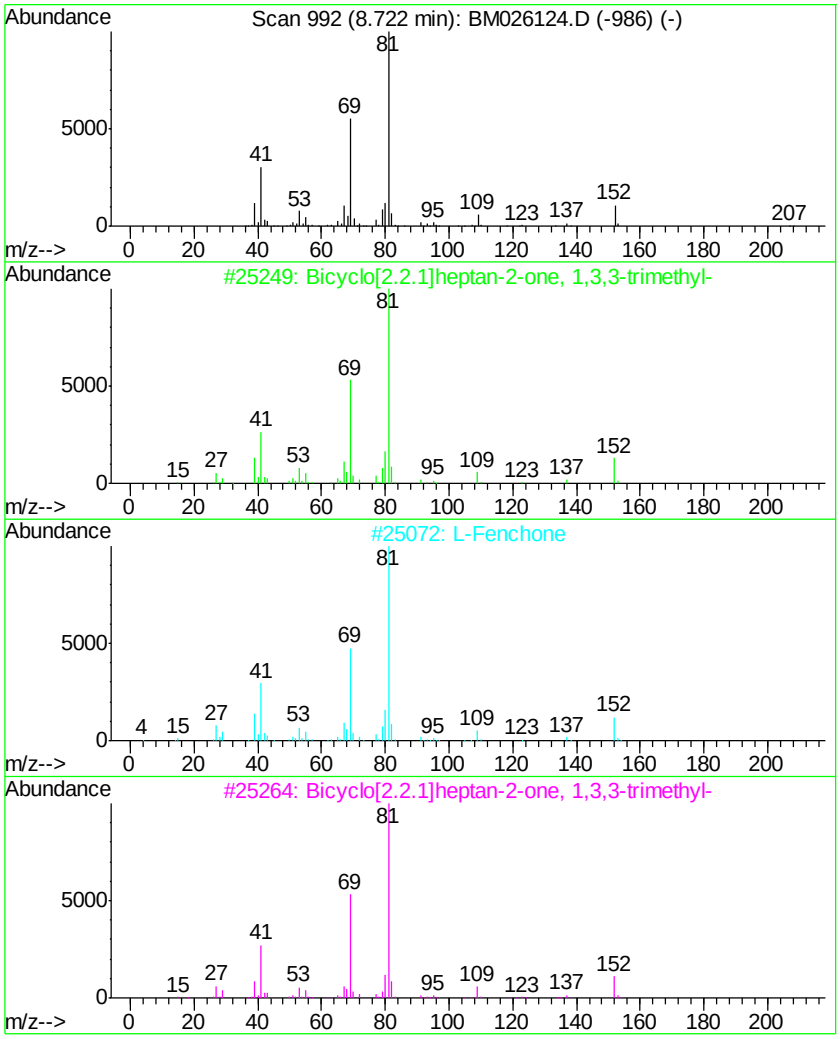
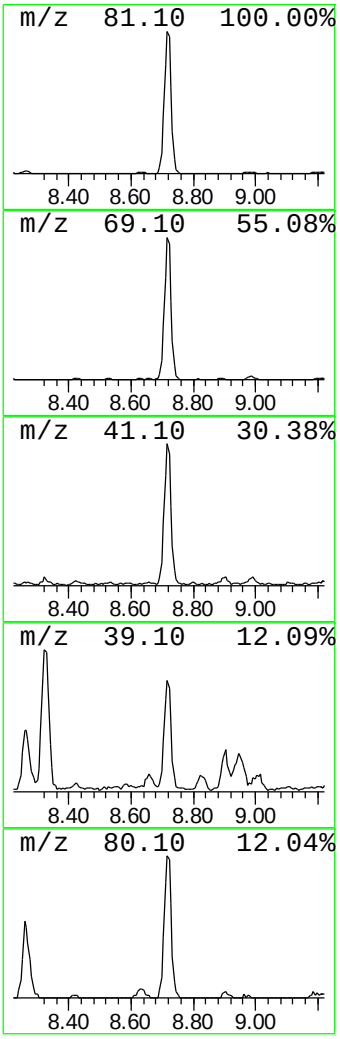
Peak Number 6 Bicyclo[2.2.1]heptan-2-one,... Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.72	6.96 ng/ul	395233	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	L-Fenchone	152	C10H16O	007787-20-4	94
3	Bicyclo[2.2.1]heptan-2-one, 1,3,...	152	C10H16O	001195-79-5	94
4	Bicyclo[2.2.1]heptan-2-one, 1,3,...	152	C10H16O	001195-79-5	91
5	D-Fenchone	152	C10H16O	004695-62-9	91





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Sample : L2737-12DL 10X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9B27DL

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Quant Title : SVOA CALIBRATION

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

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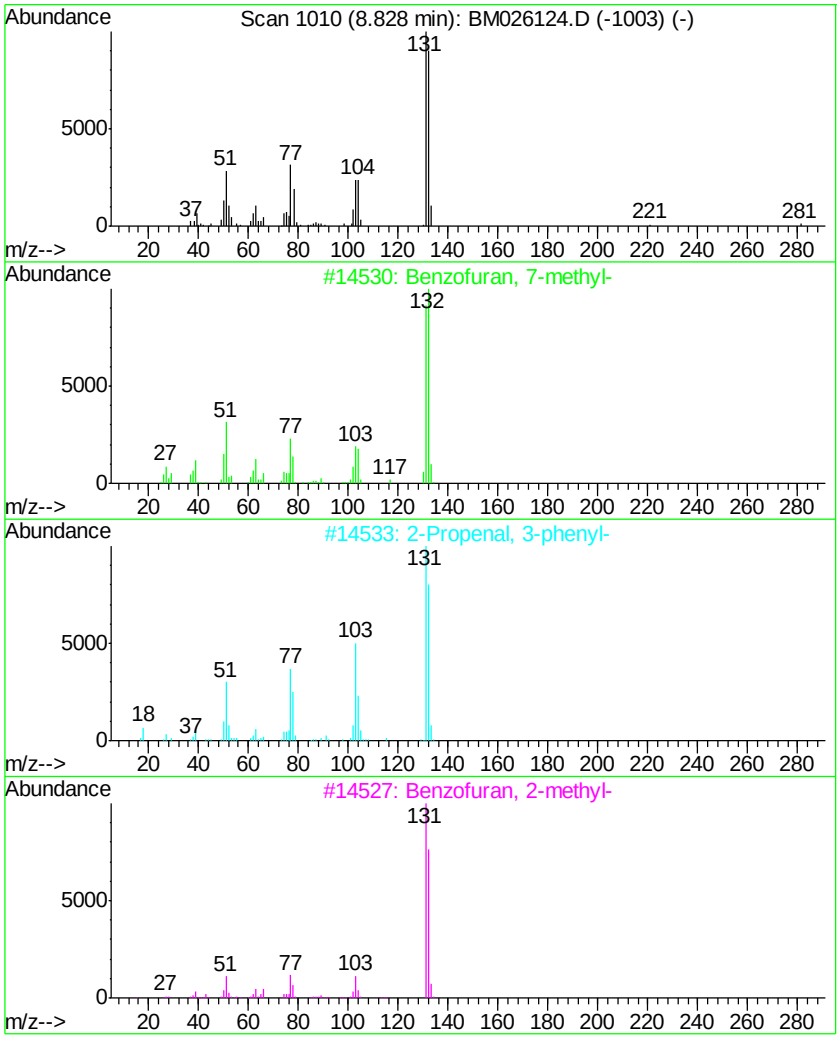
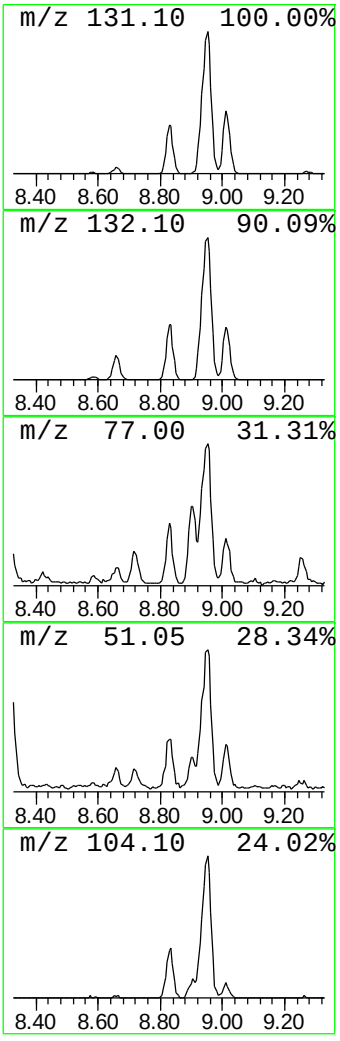
Peak Number 7 Benzofuran, 7-methyl- Concentration Rank 41

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.83	2.37 ng/ul	134680	1,4-Dichlorobenzene-d4	7.49

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzofuran, 7-methyl-	132	C9H8O	017059-52-8	94
2	2-Propenal, 3-phenyl-	132	C9H8O	000104-55-2	93
3	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	93
4	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	90
5	Cinnamaldehyde, (E)-	132	C9H8O	014371-10-9	90





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Sample : L2737-12DL 10X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9B27DL

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Quant Title : SVOA CALIBRATION

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

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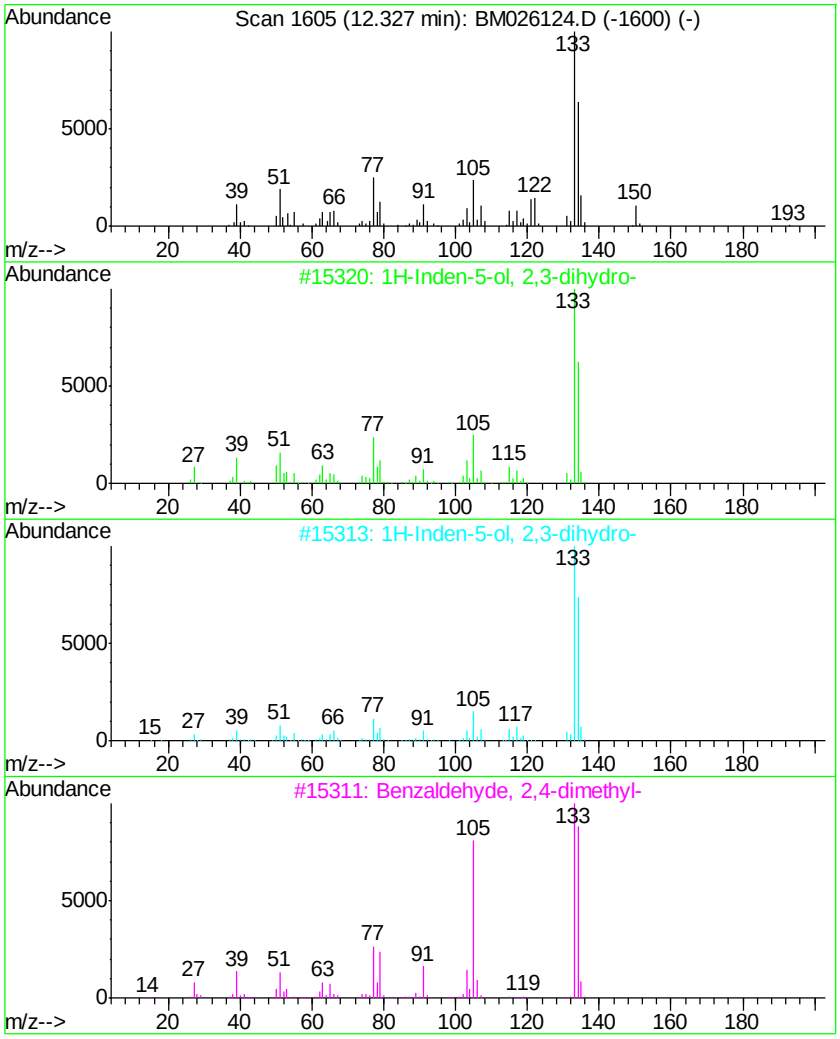
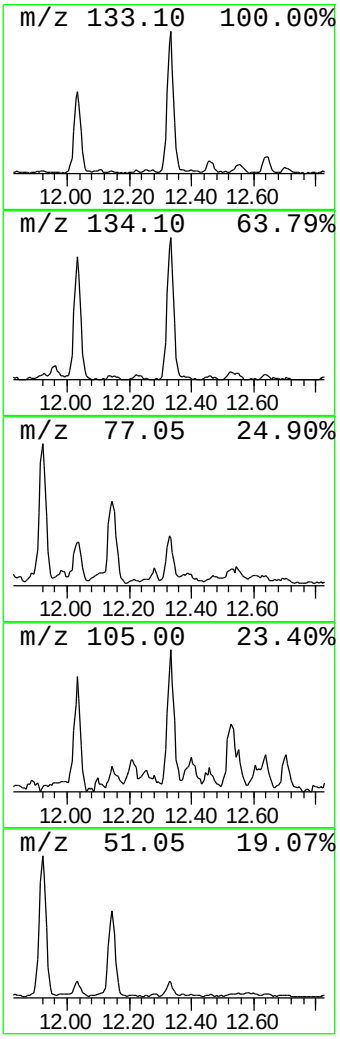
Peak Number 8 1H-Inden-5-ol, 2,3-dihydro- Concentration Rank 46

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.33	2.16 ng/ul	205326	Acenaphthene-d10	14.13

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Inden-5-ol, 2,3-dihydro-	134	C9H10O	001470-94-6	96
2	1H-Inden-5-ol, 2,3-dihydro-	134	C9H10O	001470-94-6	76
3	Benzaldehyde, 2,4-dimethyl-	134	C9H10O	015764-16-6	53
4	Phenol, 4-(2-propenyl)-	134	C9H10O	000501-92-8	53
5	1,2-Benzenediol, O-(4-ethylbenzo...	300	C17H16O5	1000330-50-2	47





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Operator : MA/SJ
Sample : L2737-12DL 10X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9B27DL

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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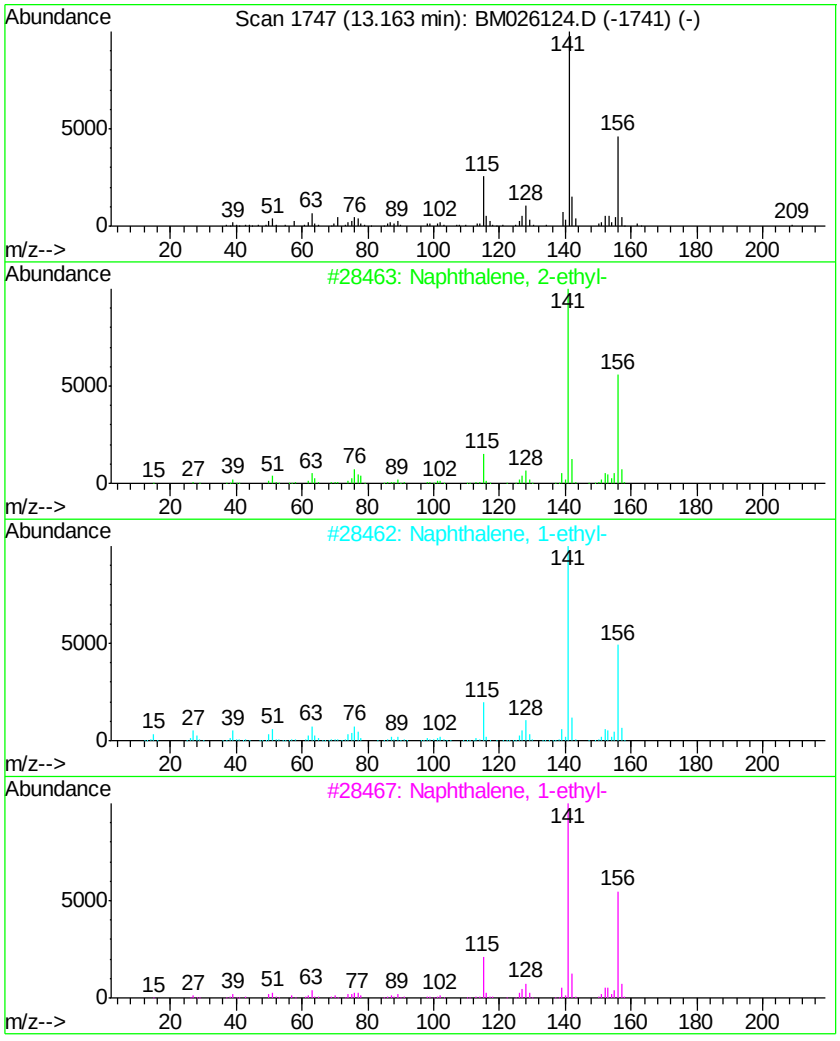
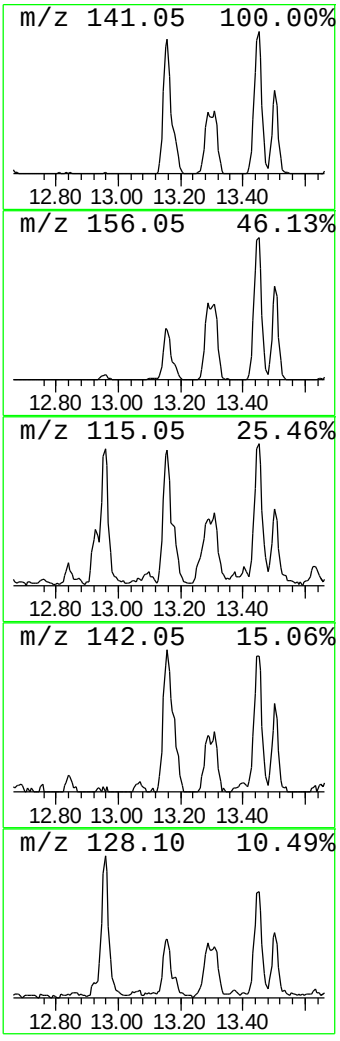
Peak Number 9 Naphthalene, 2-ethyl- Concentration Rank 32

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.16	3.48 ng/ul	331049	Acenaphthene-d10	14.13

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2-ethyl-	156	C12H12	000939-27-5	96
2	Naphthalene, 1-ethyl-	156	C12H12	001127-76-0	95
3	Naphthalene, 1-ethyl-	156	C12H12	001127-76-0	95
4	Naphthalene, 2-ethyl-	156	C12H12	000939-27-5	94
5	Naphthalene, 2-ethyl-	156	C12H12	000939-27-5	94





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Operator : MA/SJ
Sample : L2737-12DL 10X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9B27DL

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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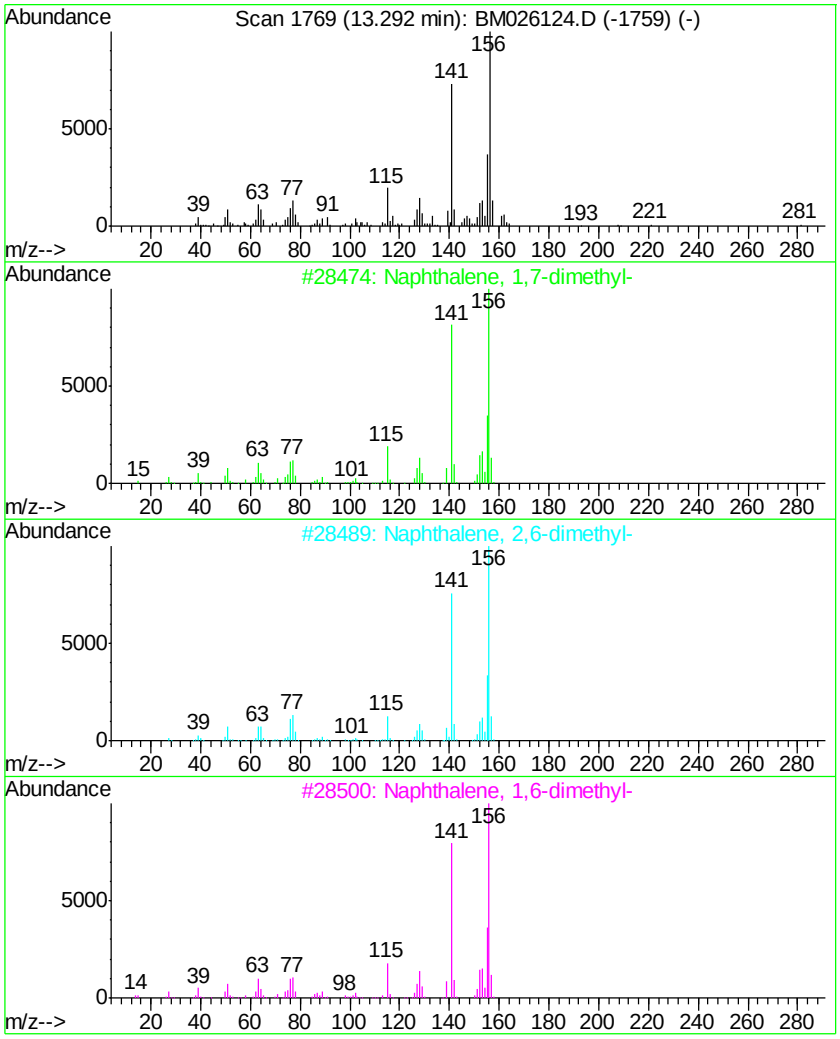
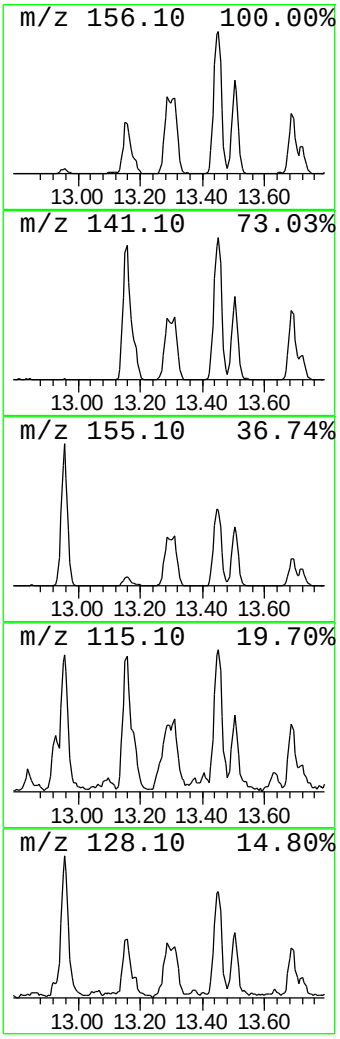
Peak Number 10 Naphthalene, 1,7-dimethyl- Concentration Rank 34

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.29	3.12 ng/ul	297130	Acenaphthene-d10	14.13

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	97
2	Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	96
3	Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	96
4	Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	96
5	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	96





Library Search Compound Report

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Sample : L2737-12DL 10X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9B27DL

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Quant Title : SVOA CALIBRATION

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

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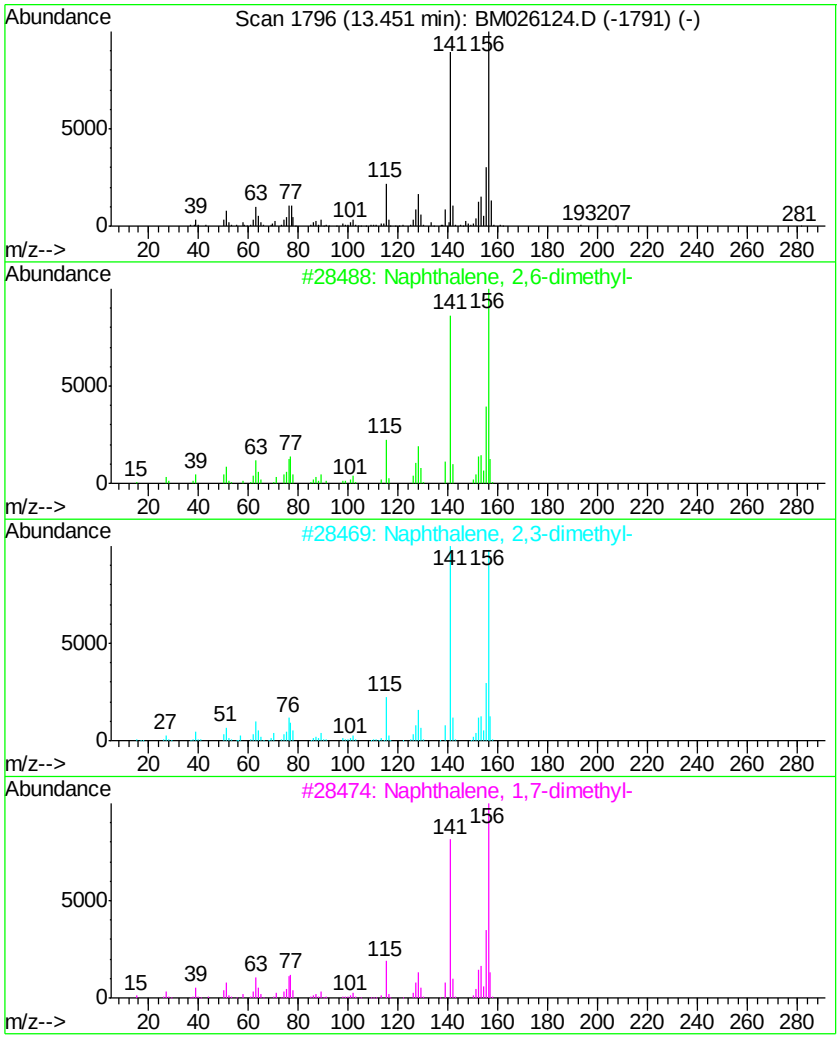
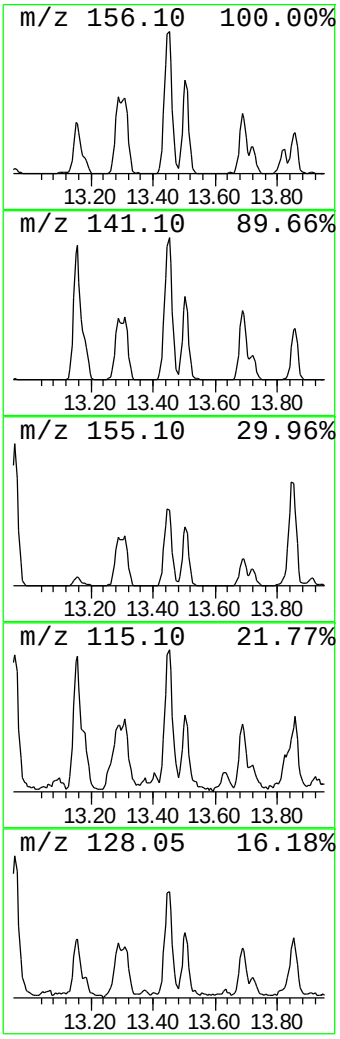
Peak Number 11 Naphthalene, 2,6-dimethyl- Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.45	4.89 ng/ul	465396	Acenaphthene-d10	14.13

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	98
2	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	98
3	Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	98
4	Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	98
5	Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	97





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Sample : L2737-12DL 10X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9B27DL

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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 12 Naphthalene, 1,5-dimethyl- Concentration Rank 38

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.50	2.52 ng/ul	239604	Acenaphthene-d10	14.13

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	98
2	Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	98
3	Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	98
4	Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	97
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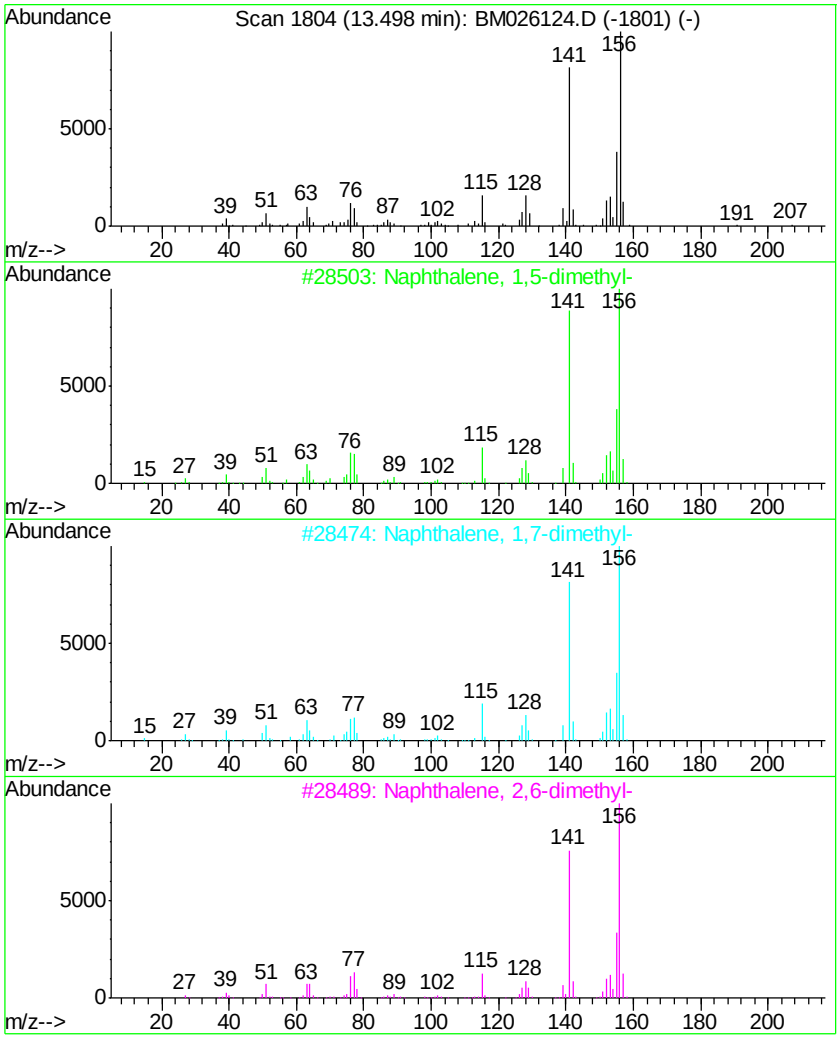

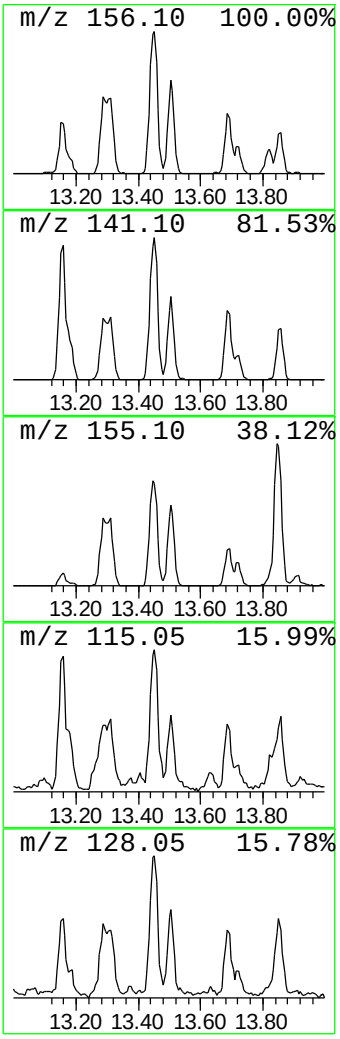

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Sample : L2737-12DL 10X
Misc :
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Instrument : BNA_M
ClientSampleId : F9B27DL

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Quant Title : SVOA CALIBRATION

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

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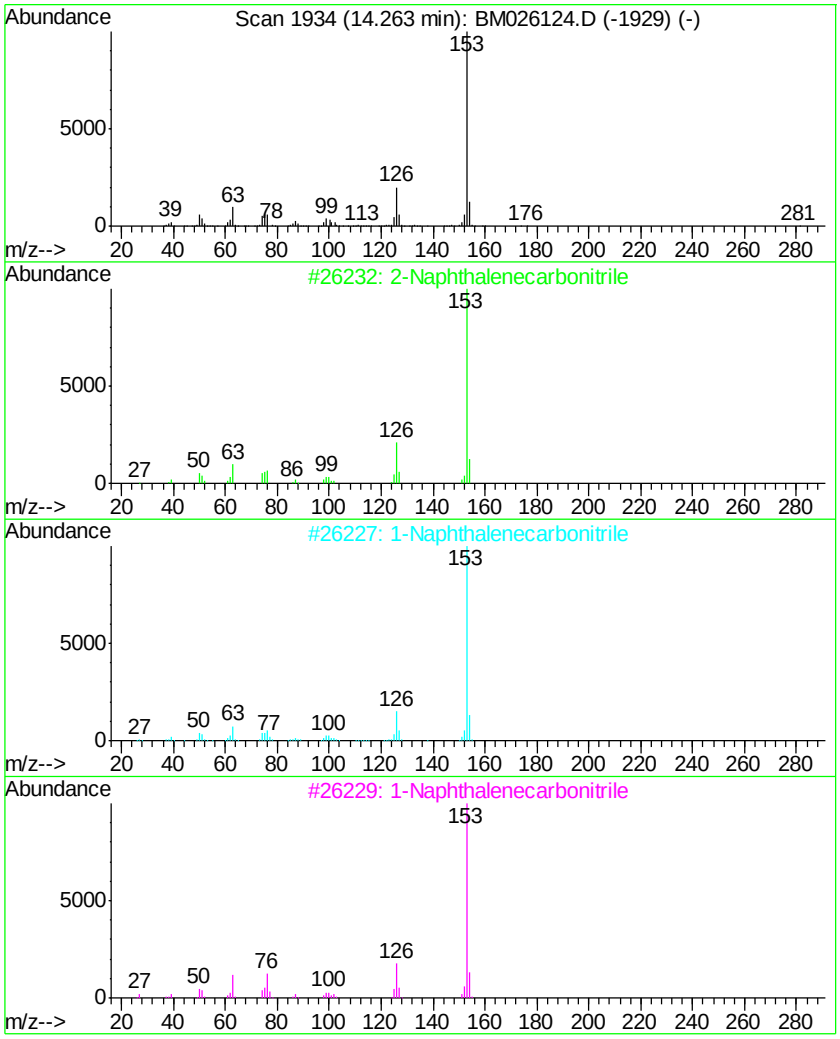
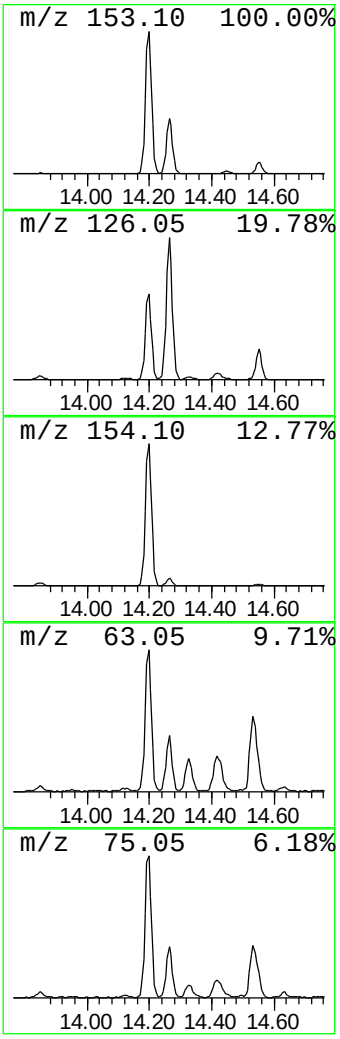
Peak Number 13 2-Naphthalenecarbonitrile Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.26	11.64 ng/ul	1108620	Acenaphthene-d10	14.13

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
4	2-Naphthalenecarbonitrile	153	C11H7N	000613-46-7	90
5	Naphthalene, 1-isocyano-	153	C11H7N	001984-04-9	90





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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 1-Naphthalenol Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.33	9.30 ng/ul	885692	Acenaphthene-d10	14.13

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Naphthalenol	144	C10H8O	000090-15-3	97
2	1-Naphthalenol	144	C10H8O	000090-15-3	94
3	1-Naphthalenol	144	C10H8O	000090-15-3	94
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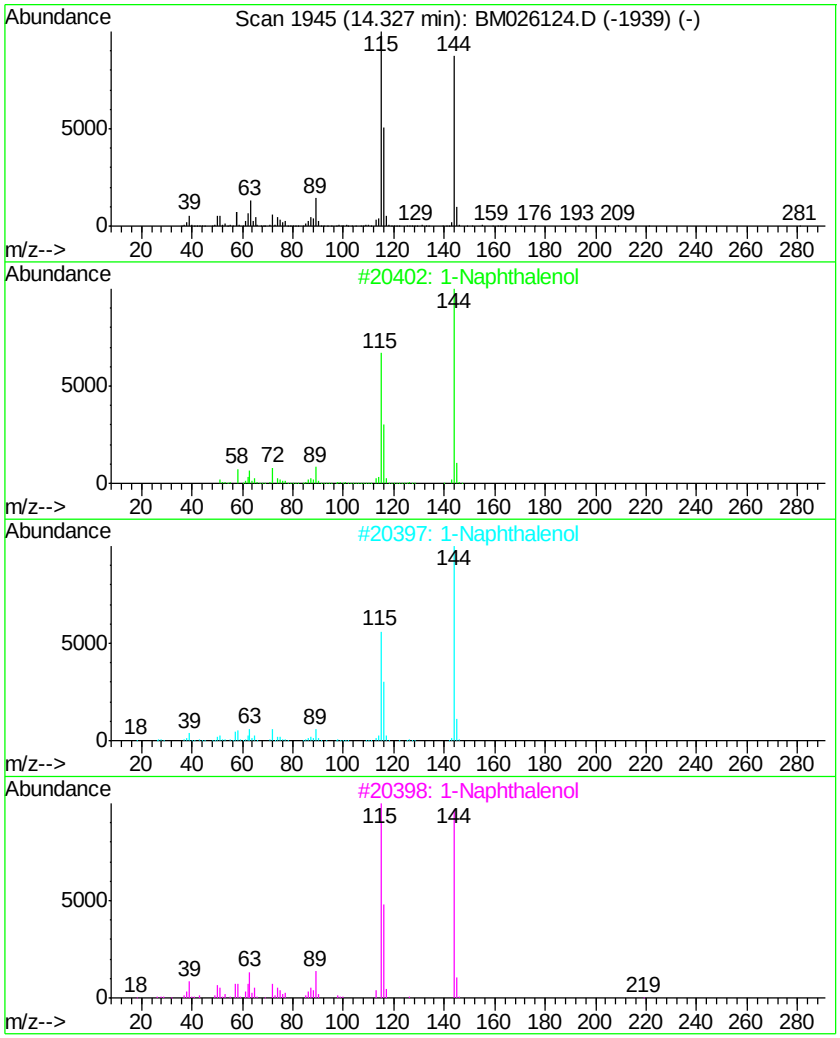
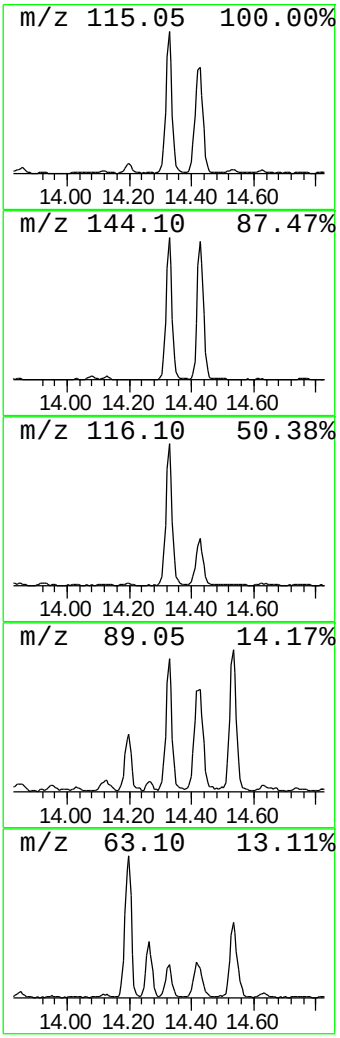
5	1-Propyne, 3-phenyl-	116	C9H8	010147-11-2	91





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Quant Title : SVOA CALIBRATION

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

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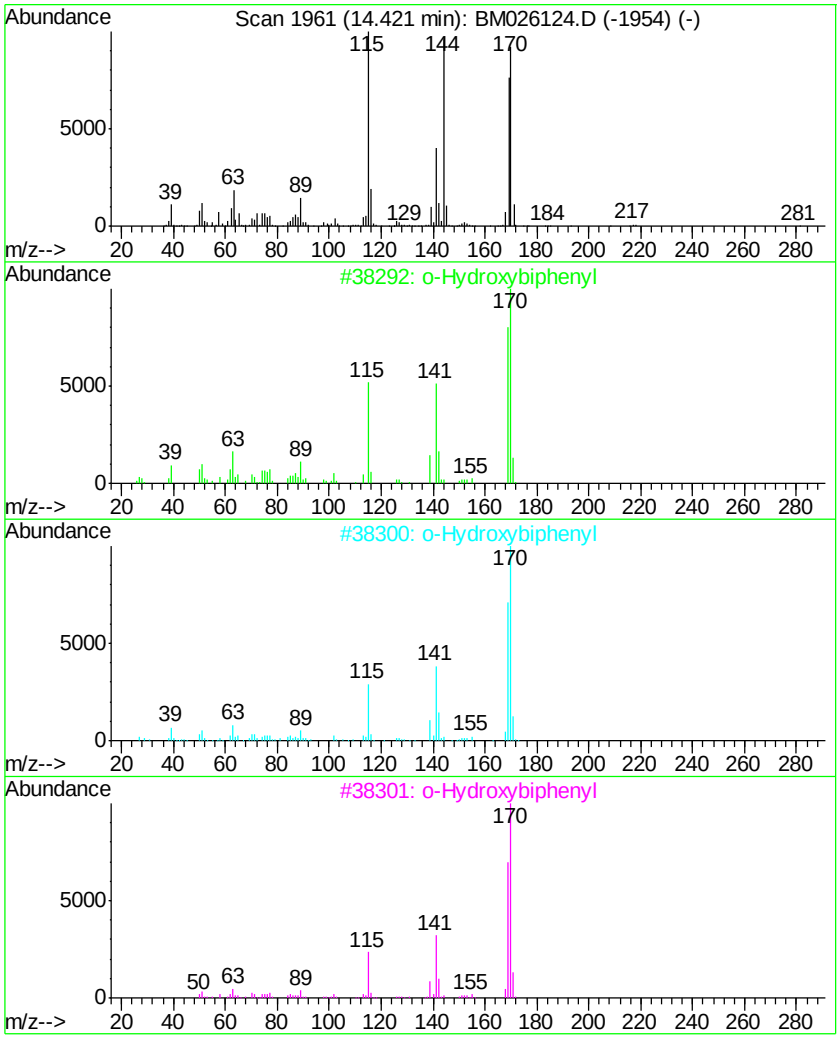
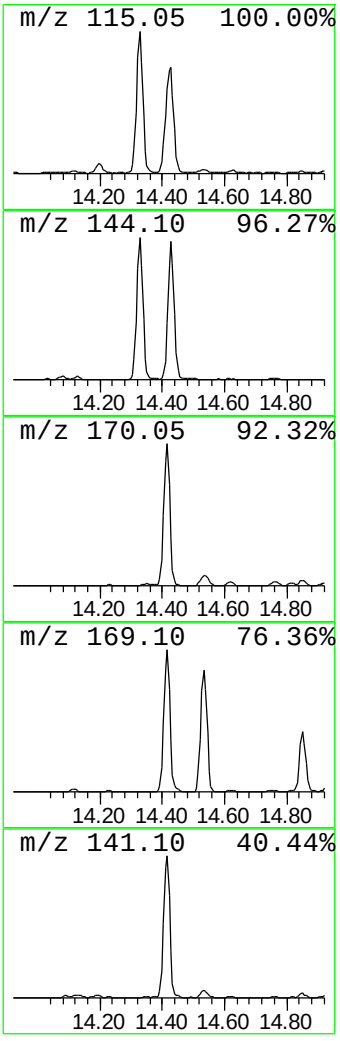
Peak Number 15 o-Hydroxybiphenyl Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.42	15.95 ng/ul	1519330	Acenaphthene-d10	14.13

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	o-Hydroxybiphenyl	170	C12H10O	000090-43-7	96
2	o-Hydroxybiphenyl	170	C12H10O	000090-43-7	93
3	o-Hydroxybiphenyl	170	C12H10O	000090-43-7	81
4	o-Hydroxybiphenyl	170	C12H10O	000090-43-7	62
5	3-Hydroxybiphenyl	170	C12H10O	000580-51-8	58





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Misc :
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Instrument : BNA_M
ClientSampleId : F9B27DL

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Quant Title : SVOA CALIBRATION

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

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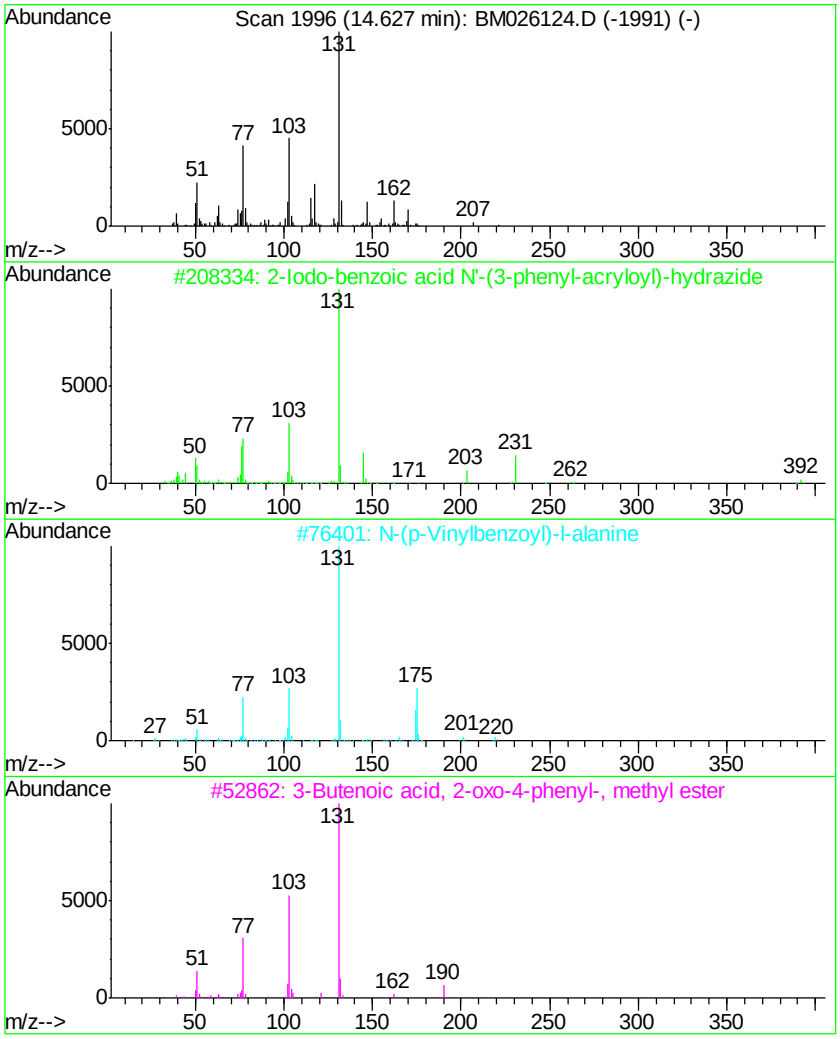
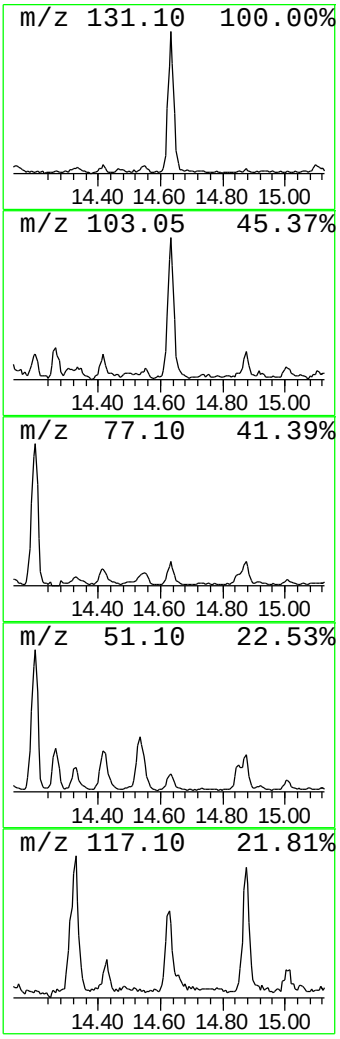
Peak Number 16 2-Iodo-benzoic acid N'-(3-p... Concentration Rank 40

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.63	2.42 ng/ul	230566	Acenaphthene-d10	14.13

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Iodo-benzoic acid N'-(3-phenyl...	392	C16H13IN2O2	315672-62-9	78
2	N-(p-Vinylbenzoyl)-l-alanine	219	C12H13NO3	139184-60-4	78
3	3-Butenoic acid, 2-oxo-4-phenyl-...	190	C11H10O3	1000142-22-4	78
4	2,2-Dimethyl-1-(2-vinylphenyl)pr...	188	C13H16O	1000210-99-9	78
5	Benzoic acid, 4-ethenyl-, methyl...	162	C10H10O2	001076-96-6	78





Library Search Compound Report

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Sample : L2737-12DL 10X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9B27DL

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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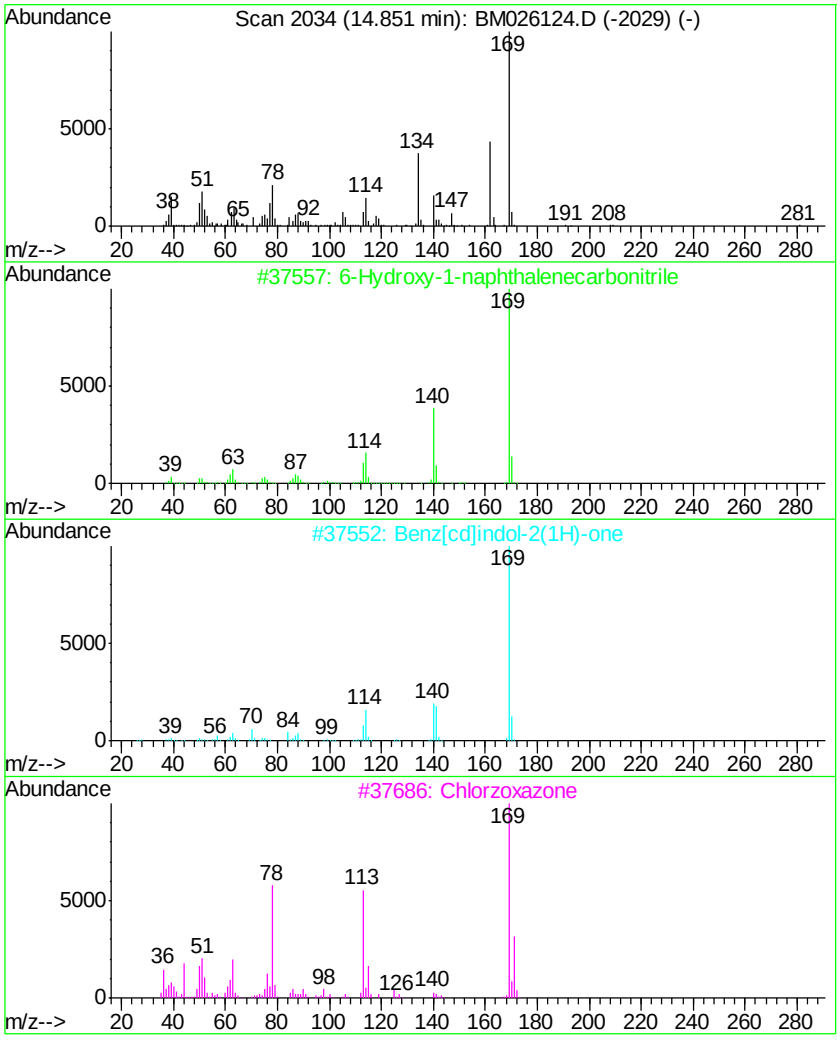
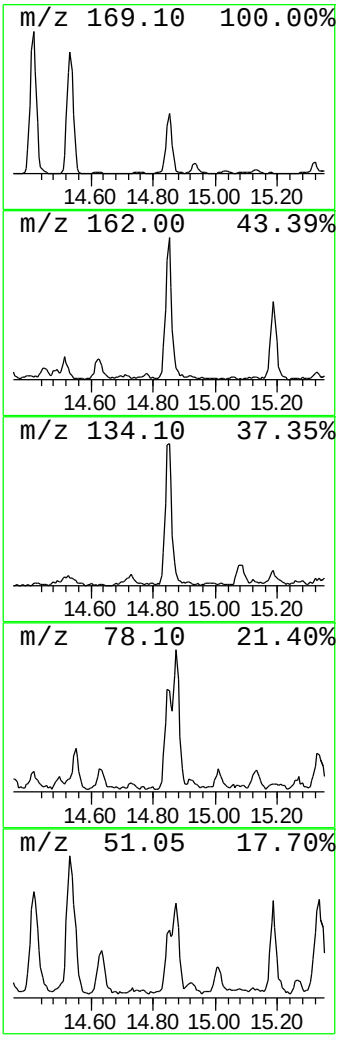
Peak Number 17 unknown-01 Concentration Rank 35

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.85	2.77 ng/ul	264010	Acenaphthene-d10	14.13

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	6-Hydroxy-1-naphthalenecarbonitrile	169	C11H7NO	1000326-74-7	43
2	Benz[cd]indol-2(1H)-one	169	C11H7NO	000130-00-7	43
3	Chlorzoxazone	169	C7H4ClNO2	000095-25-0	38
4	Naphthalene, 1-isocyanato-	169	C11H7NO	000086-84-0	38
5	Benz[cd]indol-2(1H)-one	169	C11H7NO	000130-00-7	38





Library Search Compound Report

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Data File : BM026124.D
Acq On : 26 May 2020 21:51
Operator : MA/SJ
Sample : L2737-12DL 10X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9B27DL

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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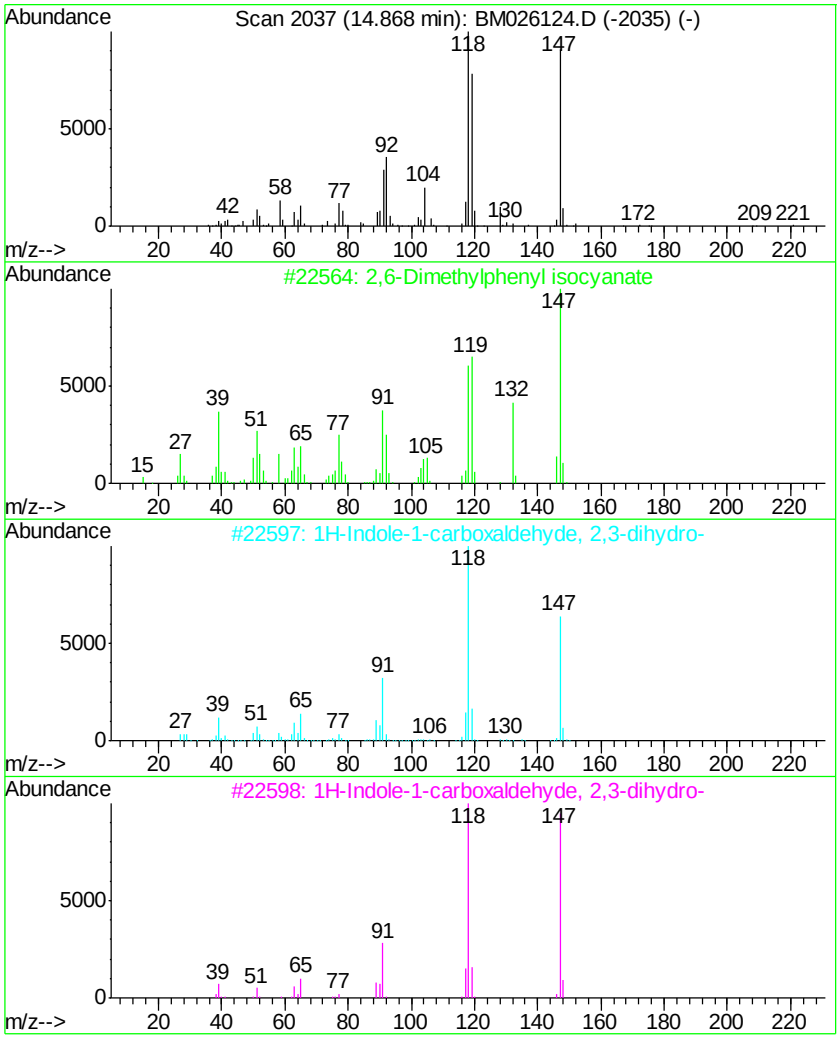
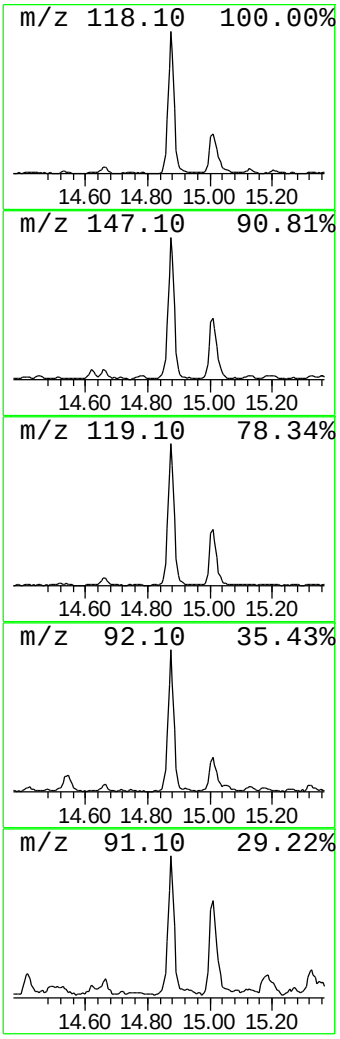
Peak Number 18 2,6-Dimethylphenyl isocyanate Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.87	5.78 ng/ul	550237	Acenaphthene-d10	14.13

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,6-Dimethylphenyl isocyanate	147	C9H9NO	028556-81-2	91
2	1H-Indole-1-carboxaldehyde, 2,3-...	147	C9H9NO	002861-59-8	70
3	1H-Indole-1-carboxaldehyde, 2,3-...	147	C9H9NO	002861-59-8	70
4	(2-Benzimidazolyl)methylamine	147	C8H9N3	005805-57-2	58
5	2-Indolinone, 1-methyl-	147	C9H9NO	000061-70-1	58





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Sample : L2737-12DL 10X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9B27DL

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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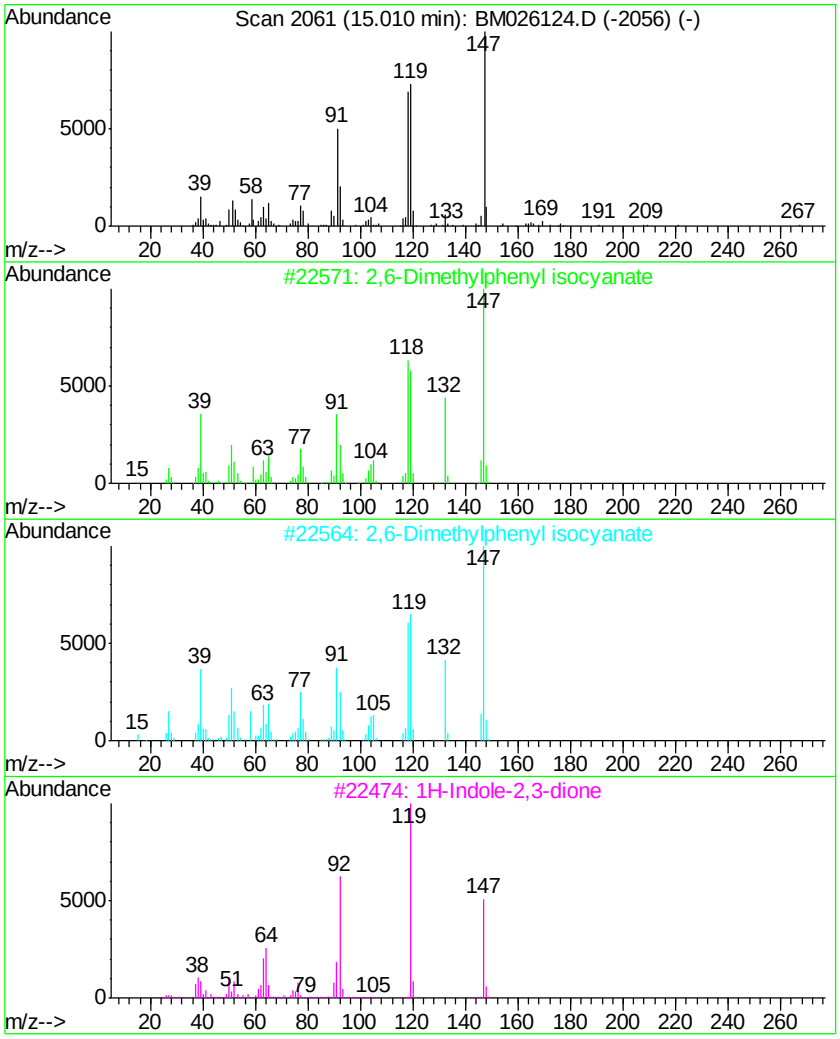
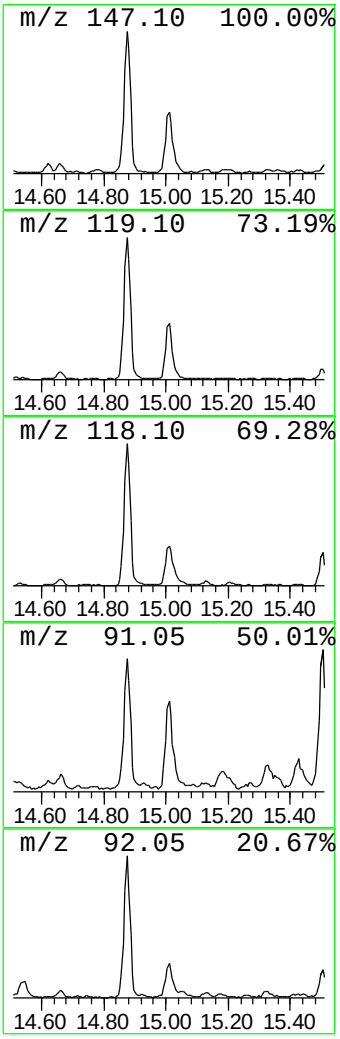
Peak Number 19 1H-Indole-2,3-dione Concentration Rank 43

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.01	2.35 ng/ul	224022	Acenaphthene-d10	14.13

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,6-Dimethylphenyl isocyanate	147	C9H9NO	028556-81-2	91
2	2,6-Dimethylphenyl isocyanate	147	C9H9NO	028556-81-2	90
3	1H-Indole-2,3-dione	147	C8H5NO2	000091-56-5	68
4	1,3,2-Benzoxazaborole, 2-ethyl-2...	147	C8H10BNO	129363-62-8	64
5	Pyrido[3,2-d]pyrimidin-4-ol	147	C7H5N3O	037538-67-3	62





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Operator : MA/SJ
Sample : L2737-12DL 10X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9B27DL

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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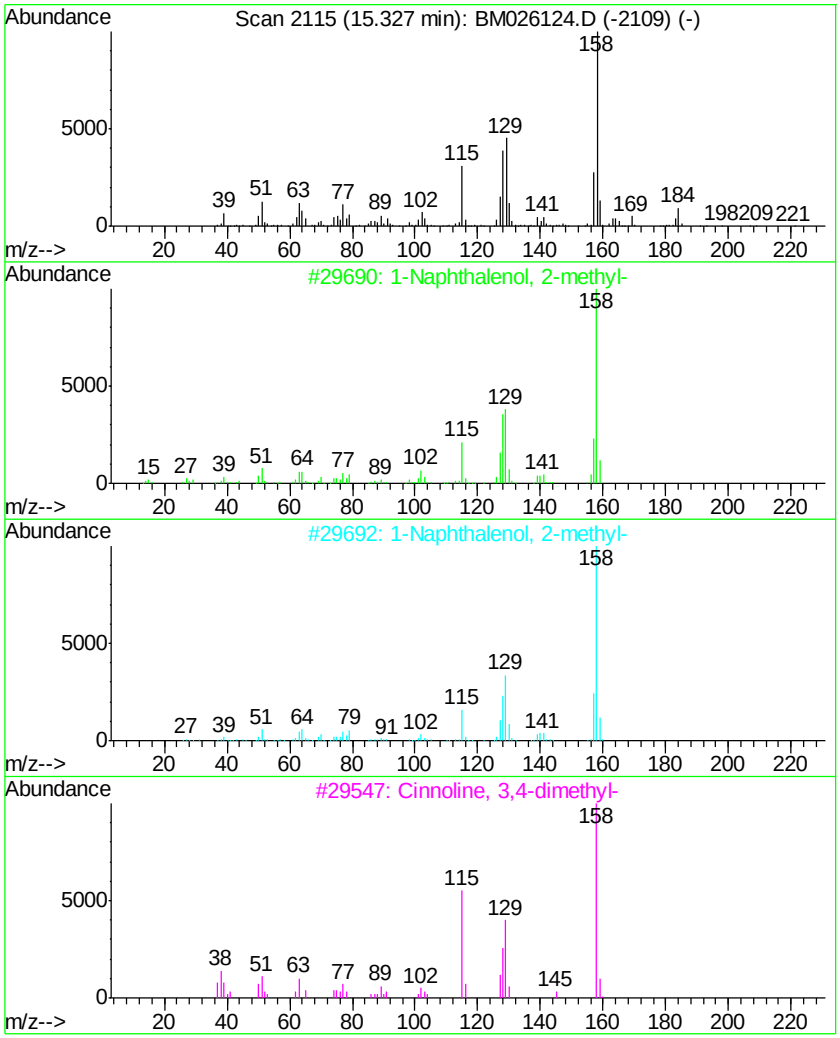
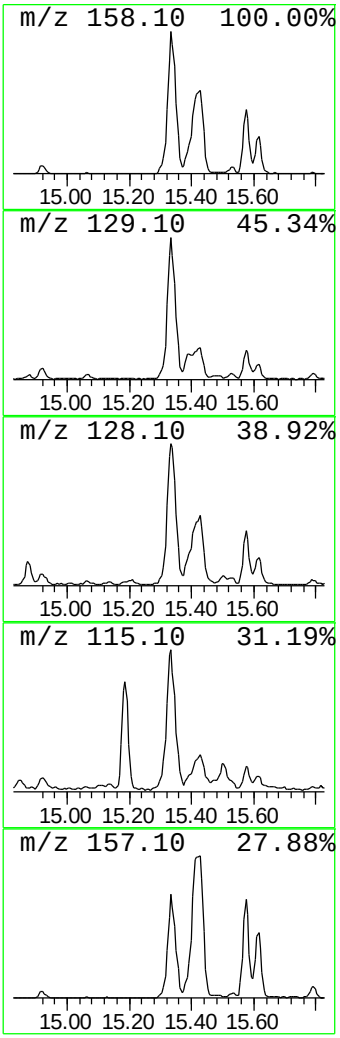
Peak Number 20 1-Naphthalenol, 2-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.33	11.80 ng/ul	1123870	Acenaphthene-d10	14.13

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Naphthalenol, 2-methyl-	158	C11H10O	007469-77-4	94
2	1-Naphthalenol, 2-methyl-	158	C11H10O	007469-77-4	93
3	Cinnoline, 3,4-dimethyl-	158	C10H10N2	003929-83-7	76
4	7-Methyl-1-naphthol	158	C11H10O	006939-33-9	58
5	1-Naphthalenol, 4-methyl-	158	C11H10O	010240-08-1	50





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Operator : MA/SJ
Sample : L2737-12DL 10X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9B27DL

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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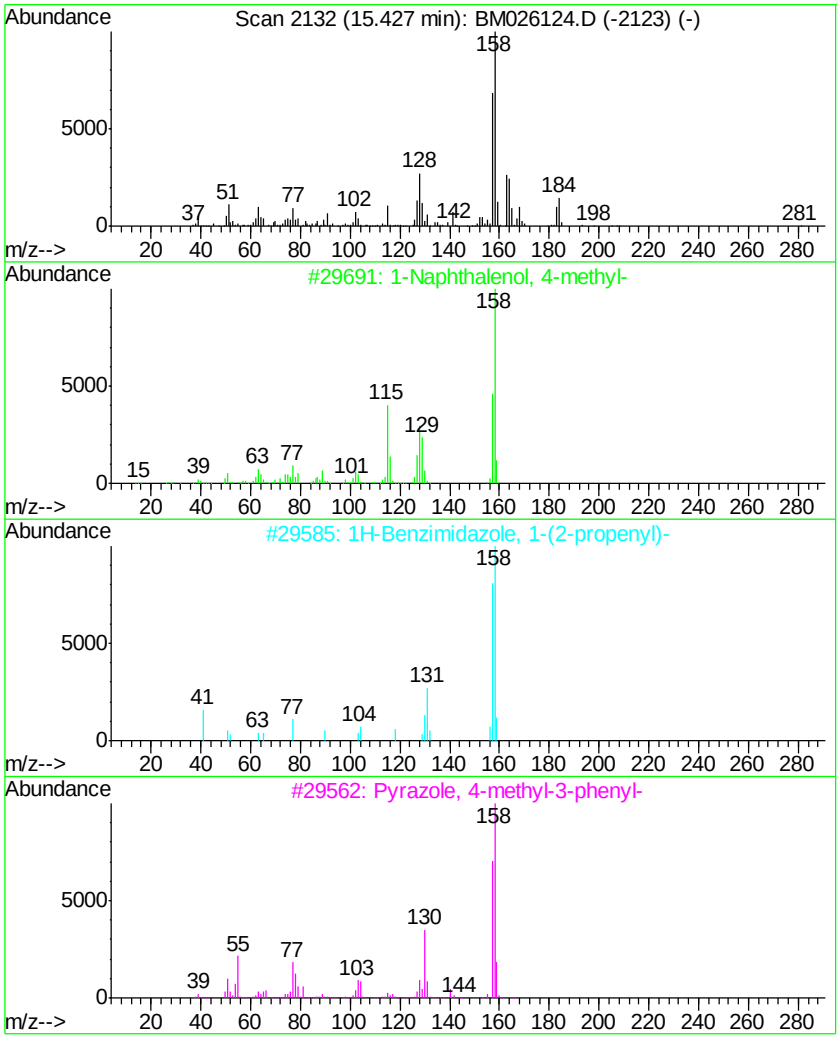
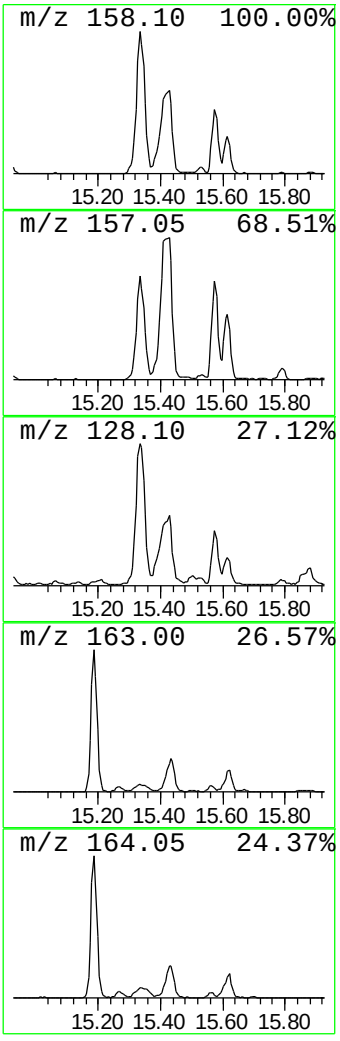
Peak Number 21 1-Naphthalenol, 4-methyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.43	9.41 ng/ul	896119	Acenaphthene-d10	14.13

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Naphthalenol, 4-methyl-	158	C11H10O	010240-08-1	50
2	1H-Benzimidazole, 1-(2-propenyl)-	158	C10H10N2	019018-22-5	43
3	Pyrazole, 4-methyl-3-phenyl-	158	C10H10N2	013808-62-3	43
4	1H-Pyrazole, 3-methyl-5-phenyl-	158	C10H10N2	003347-62-4	38
5	1-Naphthalenol, 2-methyl-	158	C11H10O	007469-77-4	38





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Operator : MA/SJ
Sample : L2737-12DL 10X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9B27DL

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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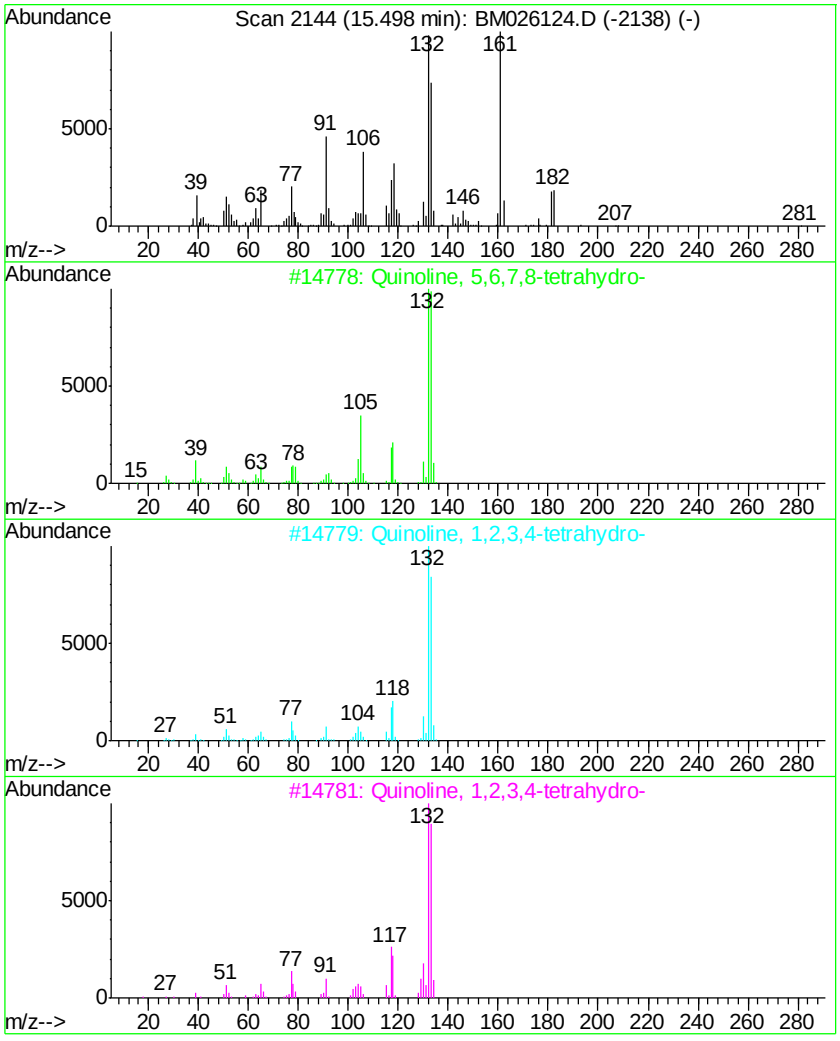
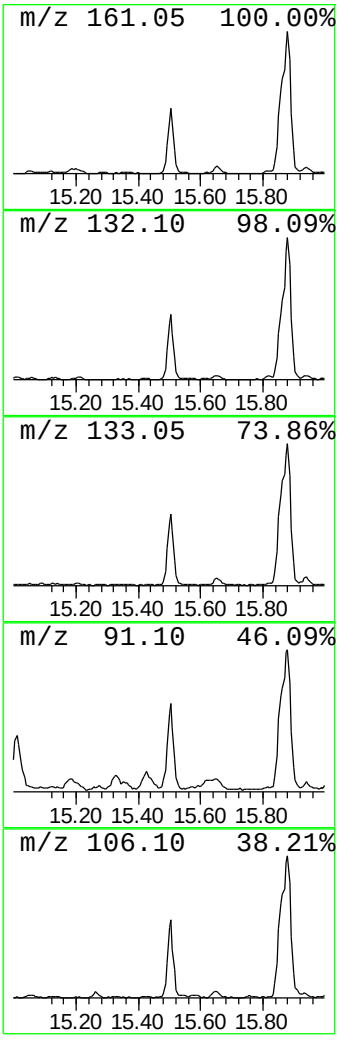
Peak Number 22 Quinoline, 5,6,7,8-tetrahydro- Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.50	6.61 ng/ul	629246	Acenaphthene-d10	14.13

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Quinoline, 5,6,7,8-tetrahydro-	133	C9H11N	010500-57-9	60
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	Mesityl isocyanate	161	C10H11NO	002958-62-5	52





Library Search Compound Report

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Sample : L2737-12DL 10X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9B27DL

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Quant Title : SVOA CALIBRATION

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

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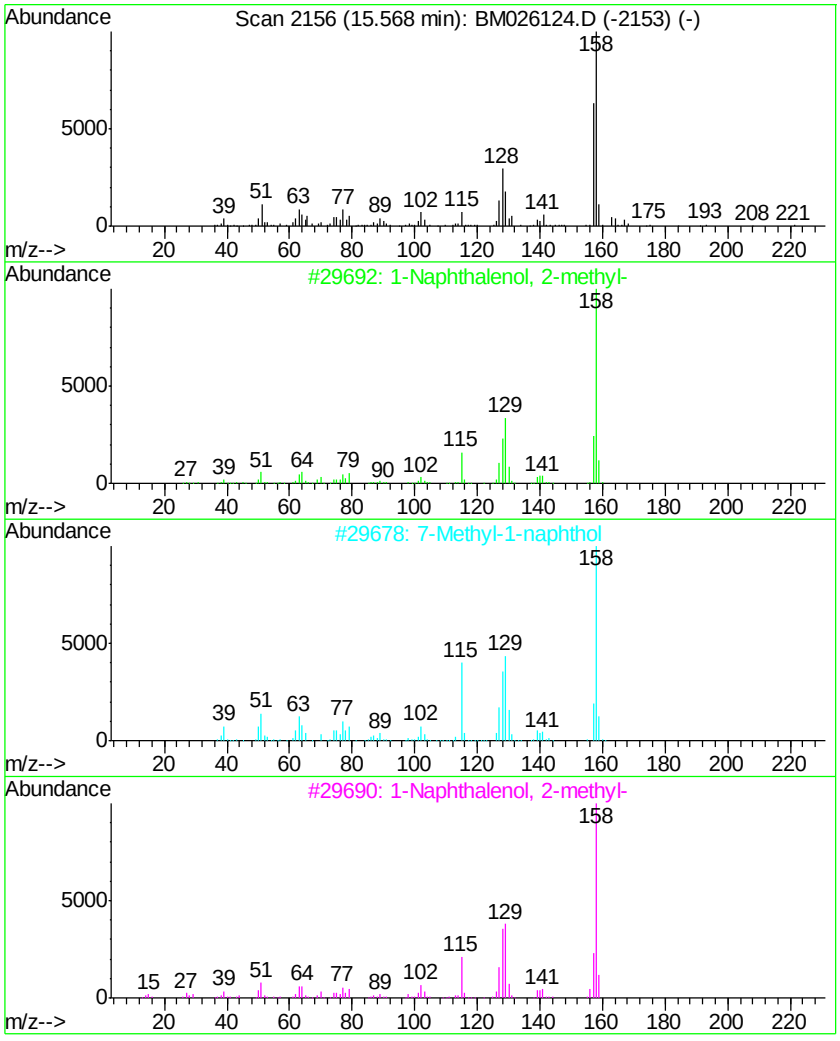
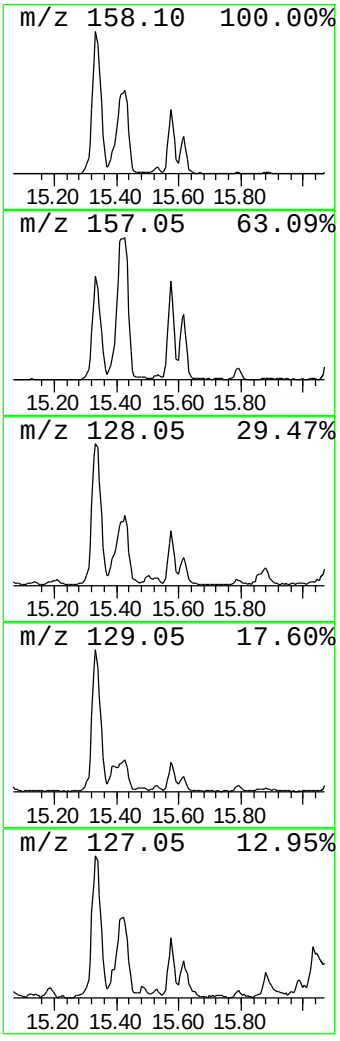
Peak Number 23 7-Methyl-1-naphthol Concentration Rank 39

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.57	2.46 ng/ul	289298	Phenanthrene-d10	16.88

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Naphthalenol, 2-methyl-	158	C11H10O	007469-77-4	59
2	7-Methyl-1-naphthol	158	C11H10O	006939-33-9	58
3	1-Naphthalenol, 2-methyl-	158	C11H10O	007469-77-4	58
4	2-Phenyl-4-methylimidazole	158	C10H10N2	000827-43-0	52
5	1H-Benzimidazole, 1-(2-propenyl)-	158	C10H10N2	019018-22-5	50





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Sample : L2737-12DL 10X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9B27DL

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Quant Title : SVOA CALIBRATION

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

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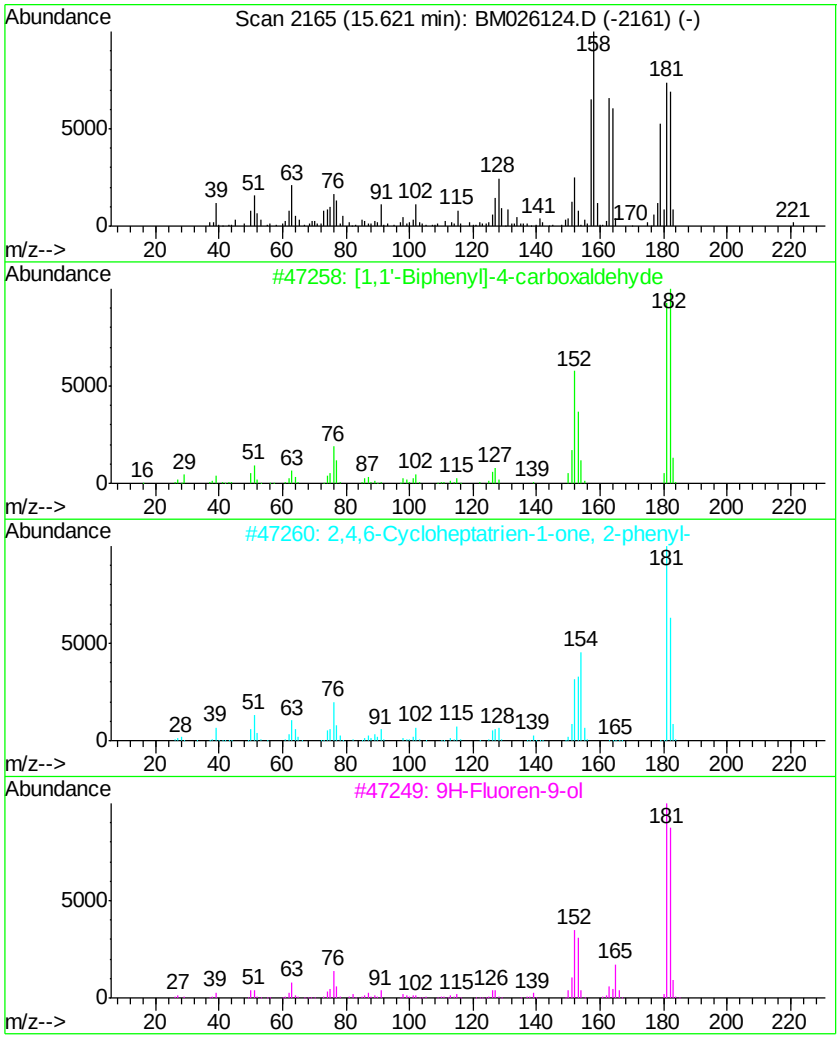
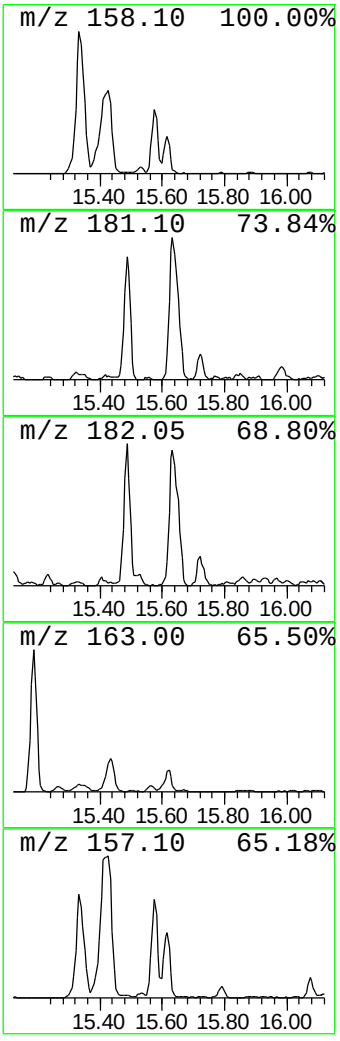
Peak Number 24 unknown-02 Concentration Rank 31

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.62	3.67 ng/ul	431643	Phenanthrene-d10	16.88

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	[1,1'-Biphenyl]-4-carboxaldehyde	182	C13H10O	003218-36-8	30
2	2,4,6-Cycloheptatrien-1-one, 2-p...	182	C13H10O	014562-09-5	30
3	9H-Fluoren-9-ol	182	C13H10O	001689-64-1	22
4	[1,1'-Biphenyl]-4-carboxaldehyde	182	C13H10O	003218-36-8	18
5	1-(2-Aminophenyl)pyrrole	158	C10H10N2	006025-60-1	18





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Sample : L2737-12DL 10X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9B27DL

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Quant Title : SVOA CALIBRATION

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

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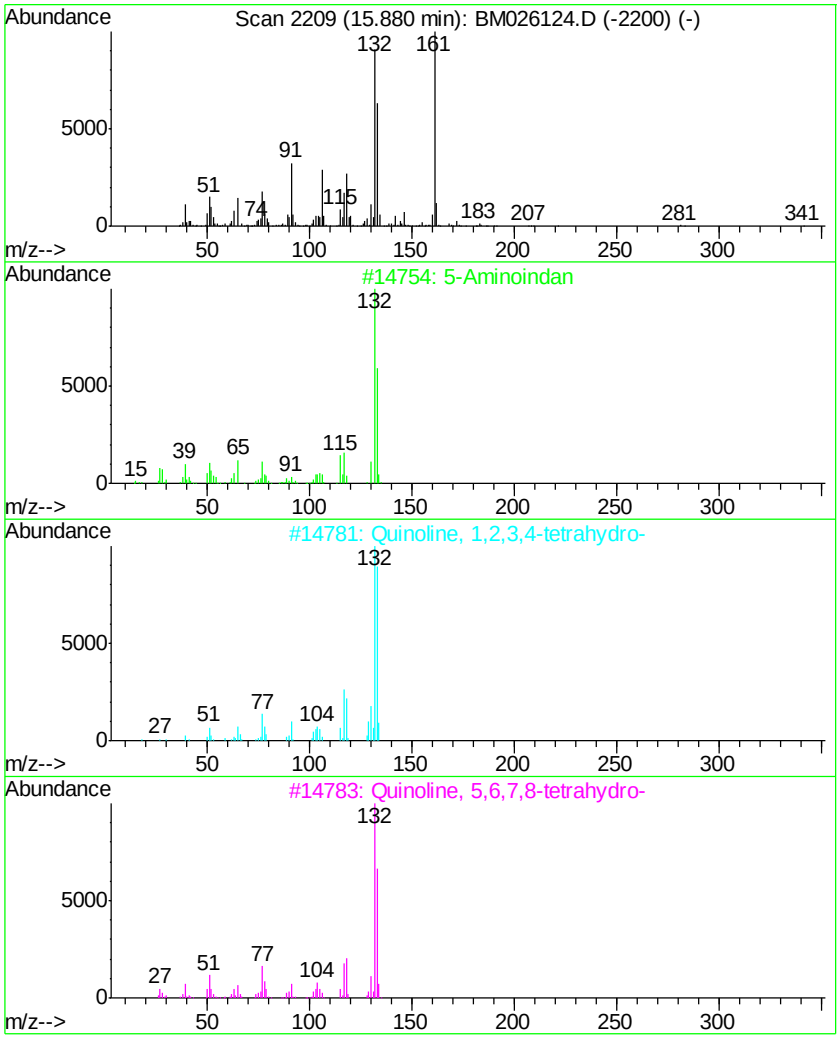
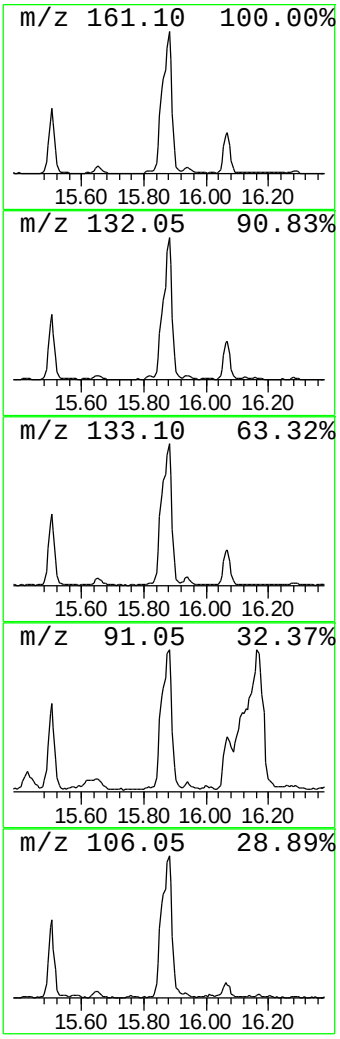
Peak Number 25 5-Aminoindan Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.88	12.12 ng/ul	1426440	Phenanthrene-d10	16.88

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5-Aminoindan	133	C9H11N	024425-40-9	55
2	Quinoline, 1,2,3,4-tetrahydro-	133	C9H11N	000635-46-1	55
3	Quinoline, 5,6,7,8-tetrahydro-	133	C9H11N	010500-57-9	55
4	Quinoline, 1,2,3,4-tetrahydro-	133	C9H11N	000635-46-1	55
5	5,6,7,8-Tetrahydroisoquinoline	133	C9H11N	1000379-32-5	53





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Sample : L2737-12DL 10X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9B27DL

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Quant Title : SVOA CALIBRATION

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

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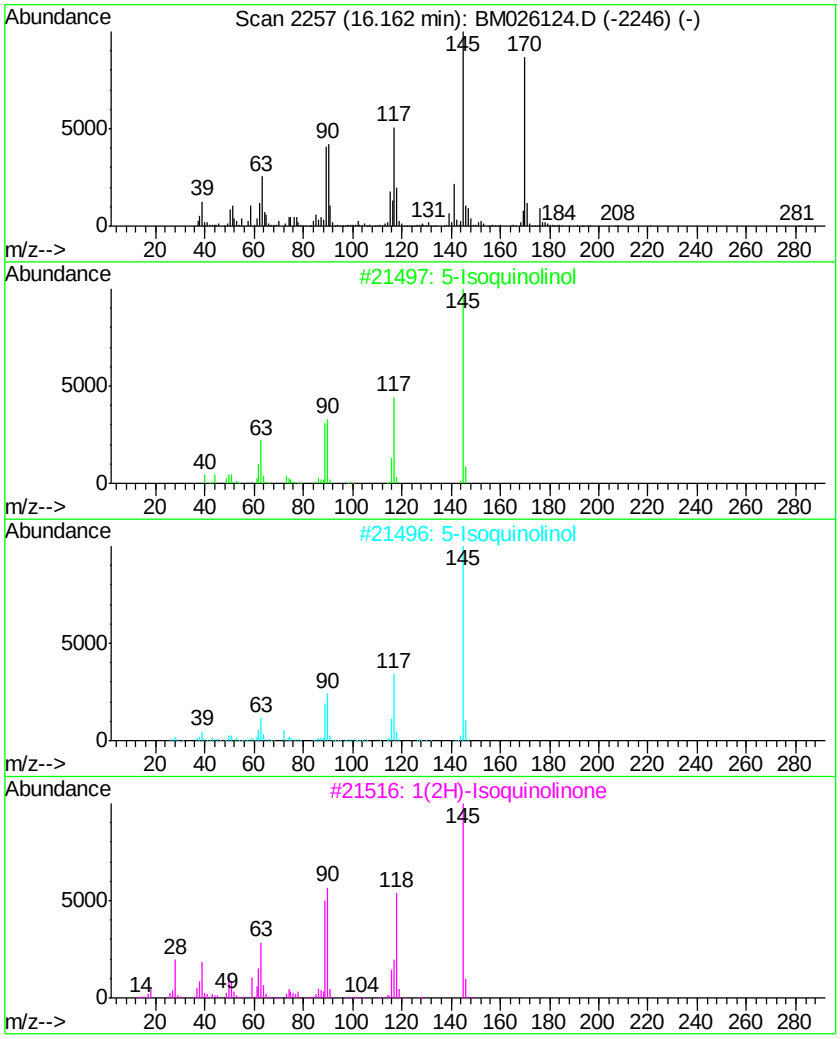
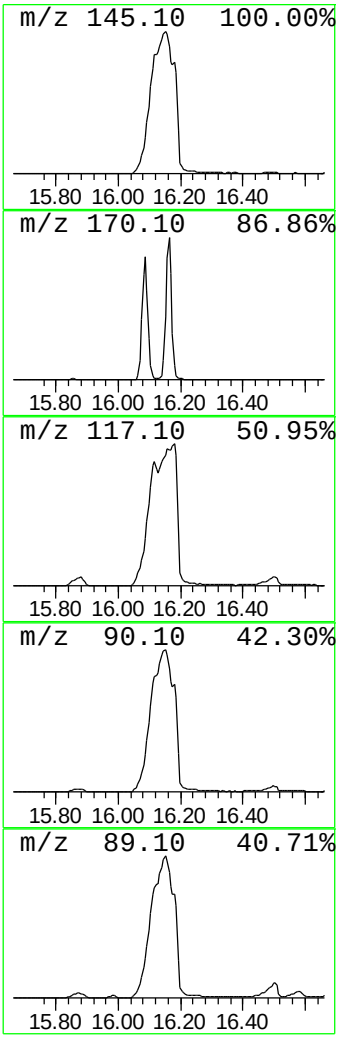
Peak Number 26 5-Isoquinolinol Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.16	83.56 ng/ul	9838840	Phenanthrene-d10	16.88

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5-Isoquinolinol	145	C9H7NO	002439-04-5	90
2	5-Isoquinolinol	145	C9H7NO	002439-04-5	90
3	1(2H)-Isoquinolinone	145	C9H7NO	000491-30-5	86
4	8-Hydroxyquinoline	145	C9H7NO	000148-24-3	78
5	Oxazole, 4-phenyl-	145	C9H7NO	020662-89-9	76





Library Search Compound Report

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Sample : L2737-12DL 10X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9B27DL

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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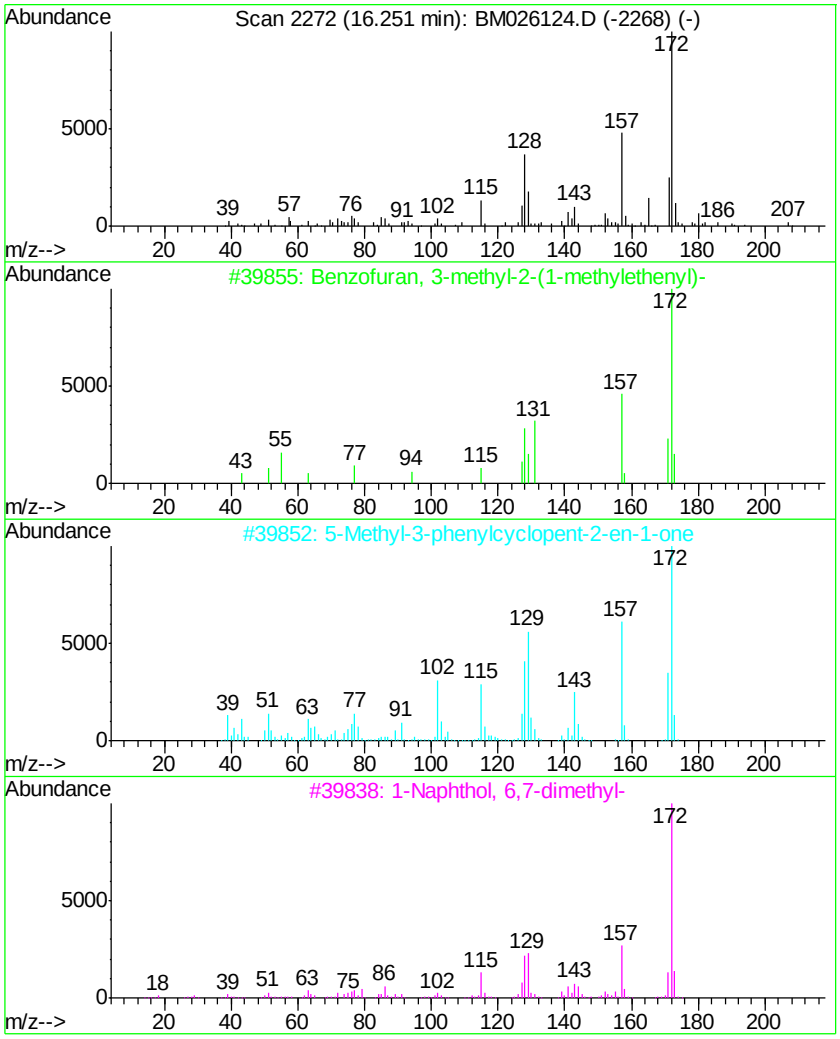
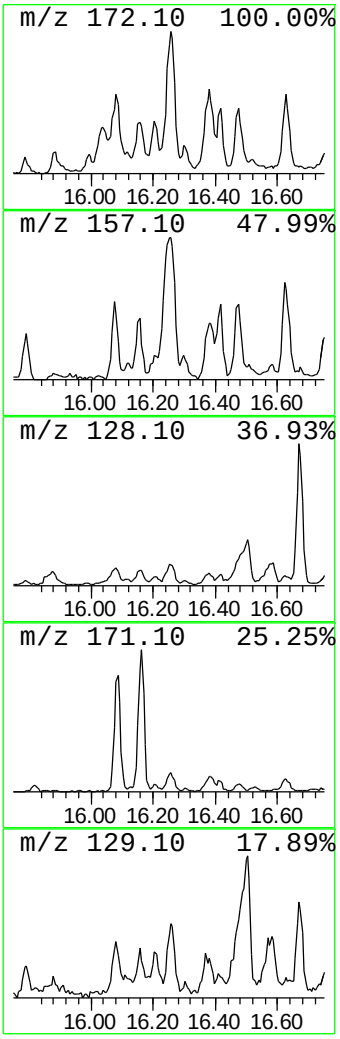
Peak Number 27 Benzofuran, 3-methyl-2-(1-m... Concentration Rank 44

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.25	2.24 ng/ul	263869	Phenanthrene-d10	16.88

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzofuran, 3-methyl-2-(1-methyl...	172	C12H12O	023911-58-2	86
2	5-Methyl-3-phenylcyclopent-2-en-...	172	C12H12O	026131-82-8	80
3	1-Naphthol, 6,7-dimethyl-	172	C12H12O	031776-14-4	59
4	1,2-Dimethoxy-3-chloro-benzene	172	C8H9ClO2	090282-99-8	52
5	Cinnoline, 4-ethyl-3-methyl-	172	C11H12N2	020873-32-9	50





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Sample : L2737-12DL 10X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument : BNA_M
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Quant Title : SVOA CALIBRATION

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

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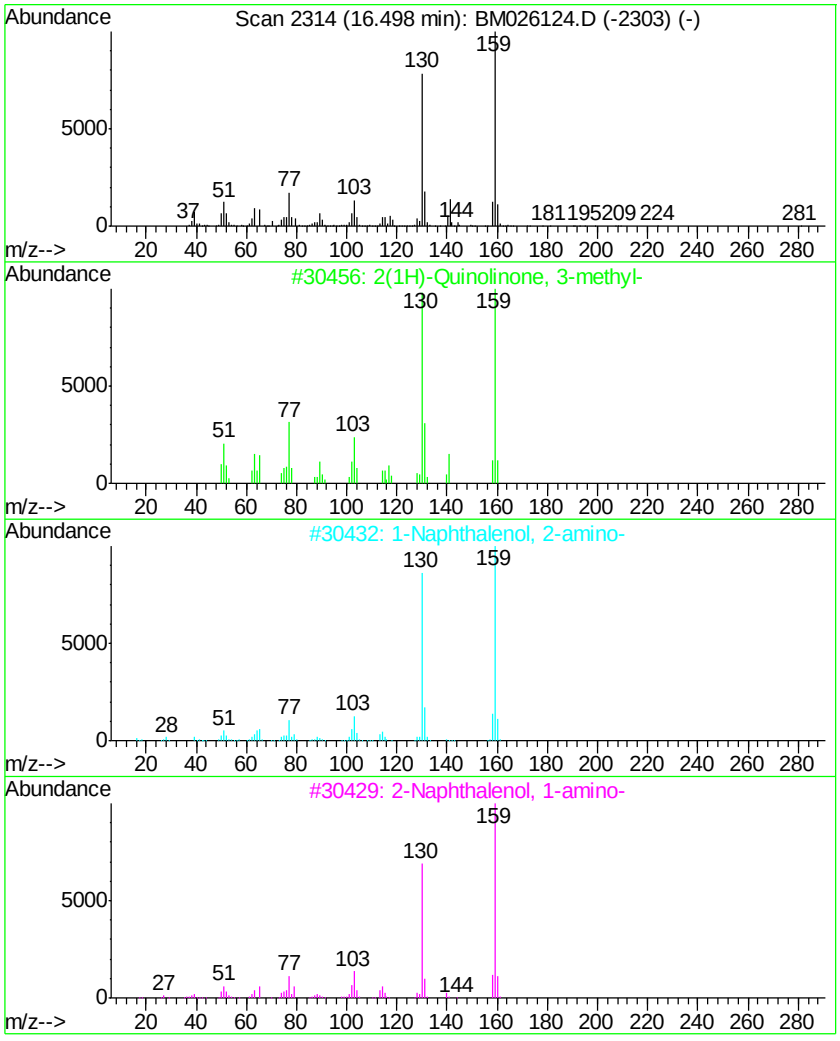
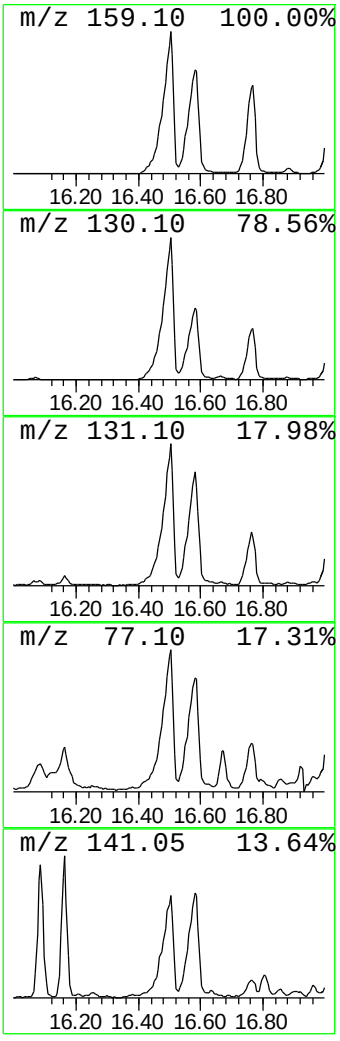
Peak Number 28 2(1H)-Quinolinone, 3-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.50	24.61 ng/ul	2897210	Phenanthrene-d10	16.88

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2(1H)-Quinolinone, 3-methyl-	159	C10H9NO	002721-59-7	97
2	1-Naphthalenol, 2-amino-	159	C10H9NO	000606-41-7	93
3	2-Naphthalenol, 1-amino-	159	C10H9NO	002834-92-6	90
4	3-Quinolinol, 2-methyl-	159	C10H9NO	000613-19-4	87
5	2(1H)-Quinolinone, 4-methyl-	159	C10H9NO	000607-66-9	87





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM052620\
Data File : BM026124.D
Acq On : 26 May 2020 21:51
Operator : MA/SJ
Sample : L2737-12DL 10X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9B27DL

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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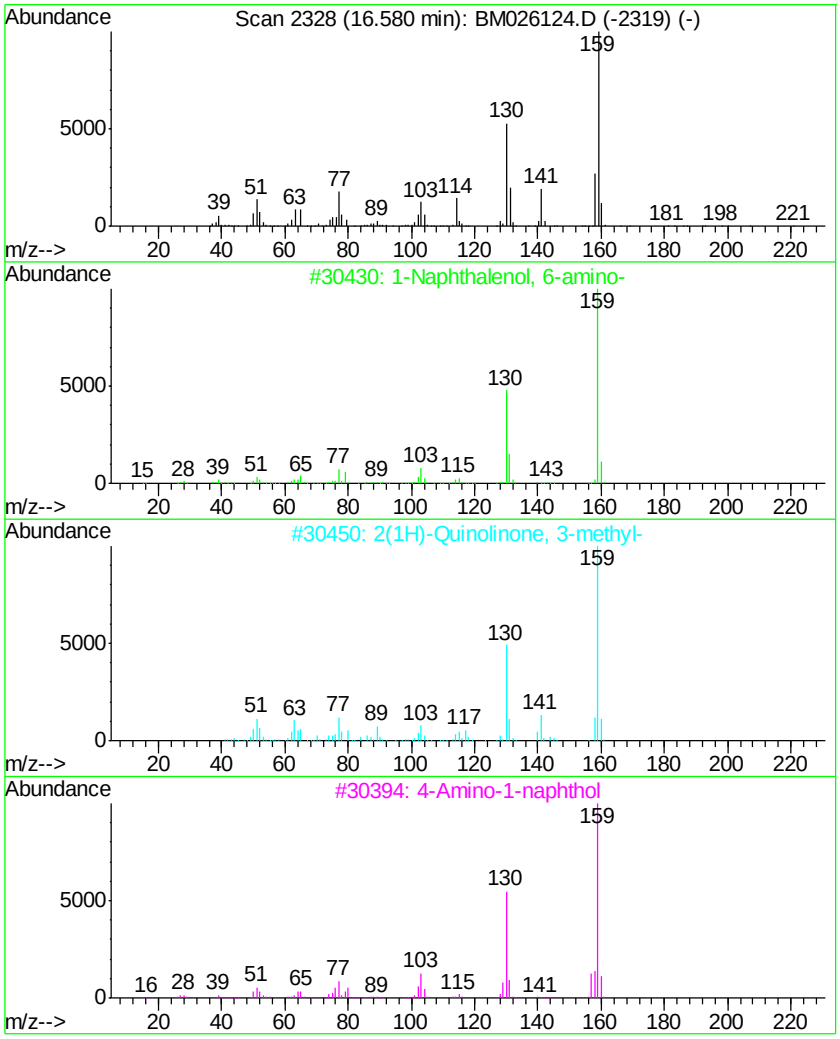
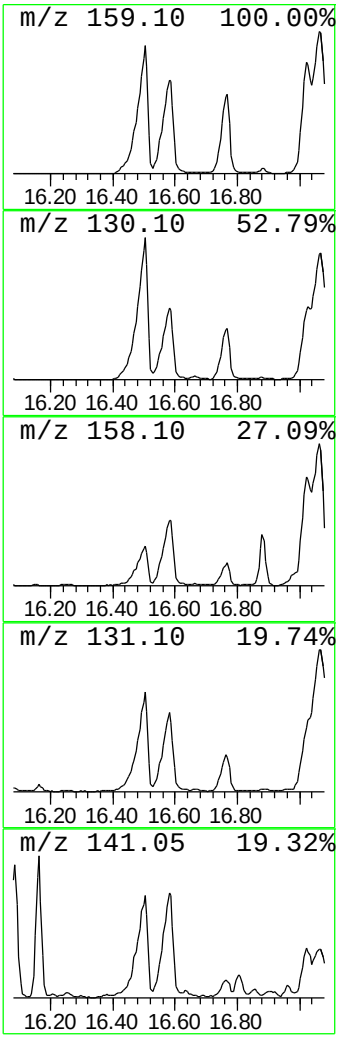
Peak Number 29 1-Naphthalenol, 6-amino- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.58	16.50 ng/ul	1942240	Phenanthrene-d10	16.88

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Naphthalenol, 6-amino-	159	C10H9NO	023894-12-4	81
2	2(1H)-Quinolinone, 3-methyl-	159	C10H9NO	002721-59-7	80
3	4-Amino-1-naphthol	159	C10H9NO	002834-90-4	76
4	2-Naphthalenol, 1-amino-	159	C10H9NO	002834-92-6	76
5	Oxazole, 4-methyl-5-phenyl-	159	C10H9NO	001008-29-3	76





Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM052620\
Data File : BM026124.D
Acq On : 26 May 2020 21:51
Operator : MA/SJ
Sample : L2737-12DL 10X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9B27DL

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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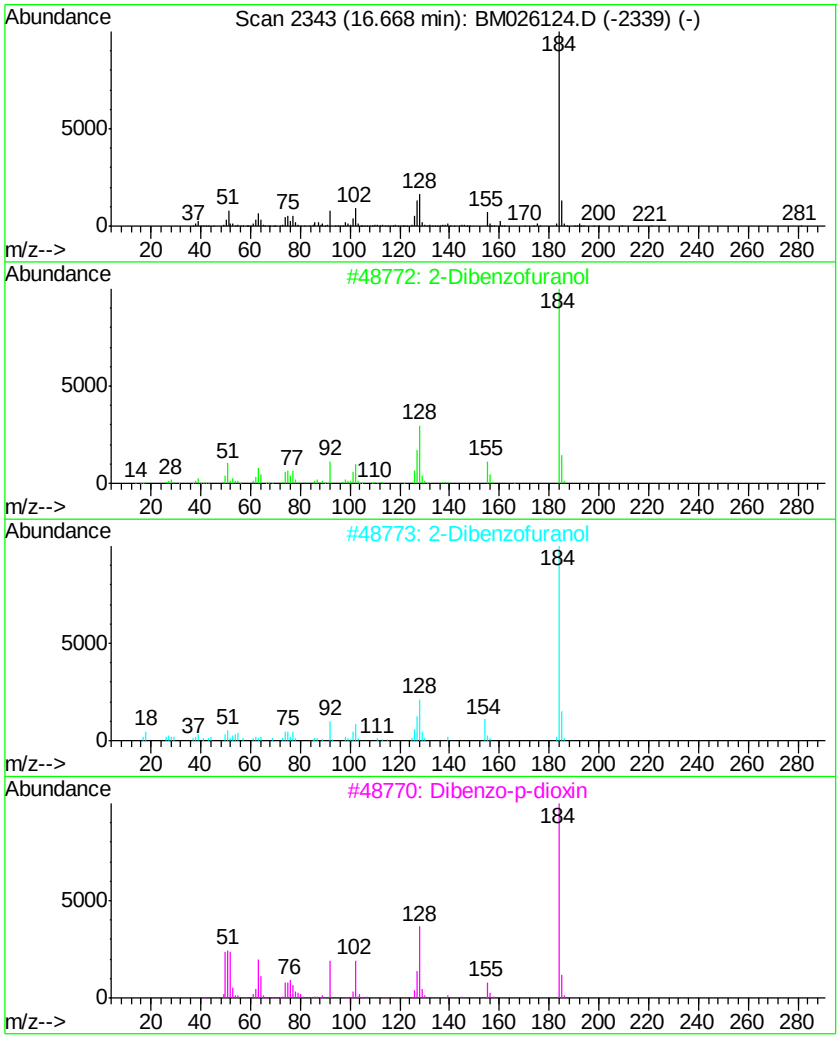
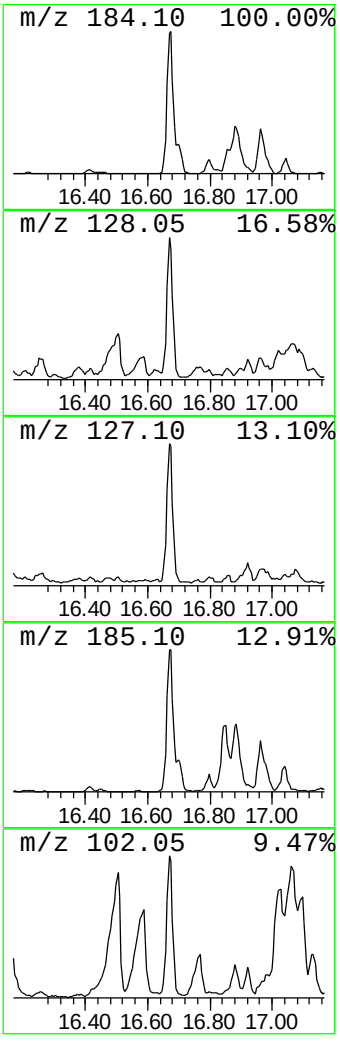
Peak Number 30 2-Dibenzofuranol Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.67	9.11 ng/ul	1073070	Phenanthrene-d10	16.88

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Dibenzofuranol	184	C12H8O2	000086-77-1	95
2	2-Dibenzofuranol	184	C12H8O2	000086-77-1	94
3	Dibenzo-p-dioxin	184	C12H8O2	000262-12-4	90
4	Dibenzo-p-dioxin	184	C12H8O2	000262-12-4	87
5	2-Dibenzofuranol	184	C12H8O2	000086-77-1	87





Tentatively Identified Compound (LSC) summary

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM052620\
Data File : BM026124.D
Acq On : 26 May 2020 21:51
Operator : MA/SJ
Sample : L2737-12DL 10X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9B27DL

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	7.21	10.9	ng/ul	621015	1	7.49	1135540	20.0
Eucalyptol	7.80	4.2	ng/ul	236472	1	7.49	1135540	20.0
Tetracyclo[3.3.1....	7.86	66.8	ng/ul	3791350	1	7.49	1135540	20.0
Indene	8.02	31.1	ng/ul	1763110	1	7.49	1135540	20.0
Benzonitrile, 4-m...	8.32	11.2	ng/ul	636743	1	7.49	1135540	20.0
Bicyclo[2.2.1]hep...	8.72	7.0	ng/ul	395233	1	7.49	1135540	20.0
Benzofuran, 7-met...	8.83	2.4	ng/ul	134680	1	7.49	1135540	20.0
1H-Inden-5-ol, 2,...	12.33	2.2	ng/ul	205326	3	14.13	1905070	20.0
Naphthalene, 2-et...	13.16	3.5	ng/ul	331049	3	14.13	1905070	20.0
Naphthalene, 1,7-...	13.29	3.1	ng/ul	297130	3	14.13	1905070	20.0
Naphthalene, 2,6-...	13.45	4.9	ng/ul	465396	3	14.13	1905070	20.0
Naphthalene, 1,5-...	13.50	2.5	ng/ul	239604	3	14.13	1905070	20.0
2-Naphthalenecarb...	14.26	11.6	ng/ul	1108620	3	14.13	1905070	20.0
1-Naphthalenol	14.33	9.3	ng/ul	885692	3	14.13	1905070	20.0
o-Hydroxybiphenyl	14.42	15.9	ng/ul	1519330	3	14.13	1905070	20.0
2-Iodo-benzoic ac...	14.63	2.4	ng/ul	230566	3	14.13	1905070	20.0
unknown-01	14.85	2.8	ng/ul	264010	3	14.13	1905070	20.0
2,6-Dimethylpheny...	14.87	5.8	ng/ul	550237	3	14.13	1905070	20.0
1H-Indole-2,3-dione	15.01	2.4	ng/ul	224022	3	14.13	1905070	20.0
1-Naphthalenol, 2...	15.33	11.8	ng/ul	1123870	3	14.13	1905070	20.0
1-Naphthalenol, 4...	15.43	9.4	ng/ul	896119	3	14.13	1905070	20.0
Quinoline, 5,6,7,...	15.50	6.6	ng/ul	629246	3	14.13	1905070	20.0
7-Methyl-1-naphthol	15.57	2.5	ng/ul	289298	4	16.88	2354820	20.0
unknown-02	15.62	3.7	ng/ul	431643	4	16.88	2354820	20.0
5-Aminoindan	15.88	12.1	ng/ul	1426440	4	16.88	2354820	20.0
5-Isoquinolinol	16.16	83.6	ng/ul	9838840	4	16.88	2354820	20.0
Benzofuran, 3-met...	16.25	2.2	ng/ul	263869	4	16.88	2354820	20.0
2(1H)-Quinolinone...	16.50	24.6	ng/ul	2897210	4	16.88	2354820	20.0
1-Naphthalenol, 6...	16.58	16.5	ng/ul	1942240	4	16.88	2354820	20.0
2-Dibenzofuranol	16.67	9.1	ng/ul	1073070	4	16.88	2354820	20.0