LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM052620\
Data File : BM026113.D
Acq On : 26 May 2020 14:26
Operator : MA/SJ
Sample : L2737-16DL 10X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9B13DL

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	6.710	641	650	665	rVB	2448965	3741438	19.41%	5.143%
2	6.840	667	672	679	rVB	53080	77223	0.40%	0.106%
3	7.028	700	704	713	rVB2	61276	96770	0.50%	0.133%
4	7.210	730	735	742	rVB	58523	92872	0.48%	0.128%
5	7.487	770	782	789	rBV	593332	902918	4.68%	1.241%
6	7.857	838	845	852	rVB	606628	925098	4.80%	1.272%
7	7.940	852	859	865	rBV	2159588	3282371	17.03%	4.512%
8	7.993	865	868	869	rVV	56369	75257	0.39%	0.103%
9	8.016	869	872	881	rVB	86860	132443	0.69%	0.182%
10	8.204	895	904	908	rBV	43068	74811	0.39%	0.103%
11	8.275	908	916	922	rVV	3192276	5666931	29.39%	7.791%
12	8.328	922	925	932	rVB	75188	116055	0.60%	0.160%
13	8.634	972	977	986	rBV2	64388	132933	0.69%	0.183%
14	8.904	1016	1023	1027	rBV	168701	272233	1.41%	0.374%
15	8.957	1027	1032	1036	rVV3	43847	86917	0.45%	0.119%
16	9.257	1075	1083	1089	rBV	65609	104037	0.54%	0.143%
17	9.351	1092	1099	1104	rBV	47408	79278	0.41%	0.109%
18	9.457	1110	1117	1120	rBV	1052298	1639250	8.50%	2.254%
19	9.487	1120	1122	1141	rVB	646747	926259	4.80%	1.273%
20	9.640	1141	1148	1152	rBV	229157	392085	2.03%	0.539%
21	9.675	1152	1154	1155	rVV2	69706	70763	0.37%	0.097%
22	9.710	1155	1160	1165	rVV	368625	897270	4.65%	1.234%
23	9.763	1165	1169	1176	rVB	1810778	2643141	13.71%	3.634%
24	9.910	1184	1194	1202	rVB3	194752	443802	2.30%	0.610%
25	10.145	1228	1234	1241	rVV	426186	646503	3.35%	0.889%
26	10.251	1245	1252	1255	rVV	730499	1222682	6.34%	1.681%
27	10.310	1255	1262	1270	rVV	11479837	19279313	100.00%	26.504%
28	10.416	1270	1280	1292	rVB	555292	926060	4.80%	1.273%
29	10.628	1308	1316	1323	rBV3	45959	108168	0.56%	0.149%
30	10.810	1342	1347	1351	rBV	80093	124125	0.64%	0.171%
31	11.051	1380	1388	1391	rBV	65873	119373	0.62%	0.164%
32	11.098	1391	1396	1405	rVB2	273102	442923	2.30%	0.609%
33	11.187	1405	1411	1417	rVB2	68978	119201	0.62%	0.164%
34	11.287	1420	1428	1433	rBV	67523	105295	0.55%	0.145%

LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM052620\
Data File : BM026113.D
Acq On : 26 May 2020 14:26
Operator : MA/SJ
Sample : L2737-16DL 10X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9B13DL

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	11.345	1433	1438	1442	rBV2	84672	136496	0.71%	0.188%
36	11.645	1482	1489	1493	rBV3	48040	87802	0.46%	0.121%
37	11.763	1499	1509	1512	rBV2	28771	62798	0.33%	0.086%
38	11.922	1527	1536	1544	rBV	723655	1155443	5.99%	1.588%
39	12.034	1544	1555	1562	rVB4	135327	315658	1.64%	0.434%
40	12.145	1562	1574	1581	rBV	678088	1072670	5.56%	1.475%
41	12.334	1601	1606	1612	rVV	150096	220943	1.15%	0.304%
42	12.469	1624	1629	1636	rBV3	38063	70966	0.37%	0.098%
43	12.957	1709	1712	1722	rVB2	74529	146424	0.76%	0.201%
44	13.157	1741	1746	1756	rVB2	56798	120081	0.62%	0.165%
45	13.286	1761	1768	1770	rBV3	67808	121710	0.63%	0.167%
46	13.451	1790	1796	1801	rVV2	104786	190428	0.99%	0.262%
47	13.504	1801	1805	1809	rVV	65540	93286	0.48%	0.128%
48	13.557	1809	1814	1819	rVV	156155	206975	1.07%	0.285%
49	13.633	1823	1827	1832	rVV2	36476	59731	0.31%	0.082%
50	13.692	1832	1837	1840	rVV	44454	73772	0.38%	0.101%
51	13.822	1853	1859	1863	rBV	156742	255786	1.33%	0.352%
52	13.851	1863	1864	1872	rVB3	48790	58060	0.30%	0.080%
53	13.951	1872	1881	1887	rBV	73826	123371	0.64%	0.170%
54	14.133	1902	1912	1917	rBV	1100869	1608925	8.35%	2.212%
55	14.198	1917	1923	1930	rVV	1386614	1932434	10.02%	2.657%
56	14.263	1930	1934	1940	rVV	82788	125938	0.65%	0.173%
57	14.333	1940	1946	1956	rVV	205067	379108	1.97%	0.521%
58	14.428	1956	1962	1970	rVV2	306342	572610	2.97%	0.787%
59	14.533	1970	1980	1991	rVV	709897	1221876	6.34%	1.680%
60	14.869	2030	2037	2042	rBV2	160822	248510	1.29%	0.342%
61	15.004	2056	2060	2067	rVV	42584	72399	0.38%	0.100%
62	15.133	2077	2082	2086	rVV	227595	316122	1.64%	0.435%
63	15.186	2086	2091	2100	rVV	684607	939564	4.87%	1.292%
64	15.263	2100	2104	2110	rVV3	48737	93640	0.49%	0.129%
65	15.339	2110	2117	2123	rVV3	104458	280362	1.45%	0.385%
66	15.433	2123	2133	2138	rVV3	79500	224587	1.16%	0.309%
67	15.504	2138	2145	2152	rVV4	65837	160216	0.83%	0.220%
68	15.580	2152	2158	2161	rVV	40145	71292	0.37%	0.098%
69	15.627	2161	2166	2176	rVB5	59572	152851	0.79%	0.210%
70	15.863	2200	2206	2216	rBV3	172560	354101	1.84%	0.487%
71	15.986	2222	2227	2229	rBV2	39189	56907	0.30%	0.078%

LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM052620\
Data File : BM026113.D
Acq On : 26 May 2020 14:26
Operator : MA/SJ
Sample : L2737-16DL 10X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9B13DL

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	16.057	2229	2239	2246	rVV2	452635	1084374	5.62%	1.491%
73	16.157	2252	2256	2261	rBV2	103198	159291	0.83%	0.219%
74	16.404	2292	2298	2306	rBV	86409	147752	0.77%	0.203%
75	16.522	2312	2318	2326	rVB4	69601	143000	0.74%	0.197%
76	16.669	2339	2343	2345	rVV	85657	111125	0.58%	0.153%
77	16.698	2345	2348	2358	rVB	100487	148441	0.77%	0.204%
78	16.804	2361	2366	2370	rBV4	35679	61182	0.32%	0.084%
79	16.880	2370	2379	2383	rBV	1414859	1964567	10.19%	2.701%
80	16.921	2383	2386	2390	rVV	825133	1068608	5.54%	1.469%
81	16.980	2390	2396	2409	rVB4	326823	755629	3.92%	1.039%
82	17.216	2429	2436	2440	rBV3	57315	121818	0.63%	0.167%
83	17.257	2440	2443	2444	rVV	65931	72915	0.38%	0.100%
84	17.286	2444	2448	2458	rVB	290413	426500	2.21%	0.586%
85	17.386	2458	2465	2471	rBV2	176241	296738	1.54%	0.408%
86	17.451	2471	2476	2482	rVB	195592	254040	1.32%	0.349%
87	17.721	2519	2522	2526	rBV	62398	83123	0.43%	0.114%
88	17.804	2531	2536	2540	rBV2	84418	137546	0.71%	0.189%
89	17.880	2546	2549	2555	rVB	117113	153136	0.79%	0.211%
90	17.951	2555	2561	2573	rBV5	68243	136688	0.71%	0.188%
91	18.045	2573	2577	2585	rBV2	37085	95936	0.50%	0.132%
92	18.945	2726	2730	2732	rBV	161186	214898	1.11%	0.295%
93	18.968	2732	2734	2739	rVV	132823	147948	0.77%	0.203%
94	19.151	2761	2765	2771	rBV2	42741	92475	0.48%	0.127%
95	19.280	2782	2787	2790	rBV	340802	446054	2.31%	0.613%
96	19.310	2790	2792	2806	rVB3	111069	213776	1.11%	0.294%
97	19.692	2854	2857	2861	rBV	46072	61607	0.32%	0.085%
98	21.080	3088	3093	3098	rBV	1916162	2271121	11.78%	3.122%
99	23.062	3427	3430	3437	rVB	211547	319402	1.66%	0.439%
100	23.198	3448	3453	3463	rVB	1305954	2201809	11.42%	3.027%

Sum of corrected areas: 72741139

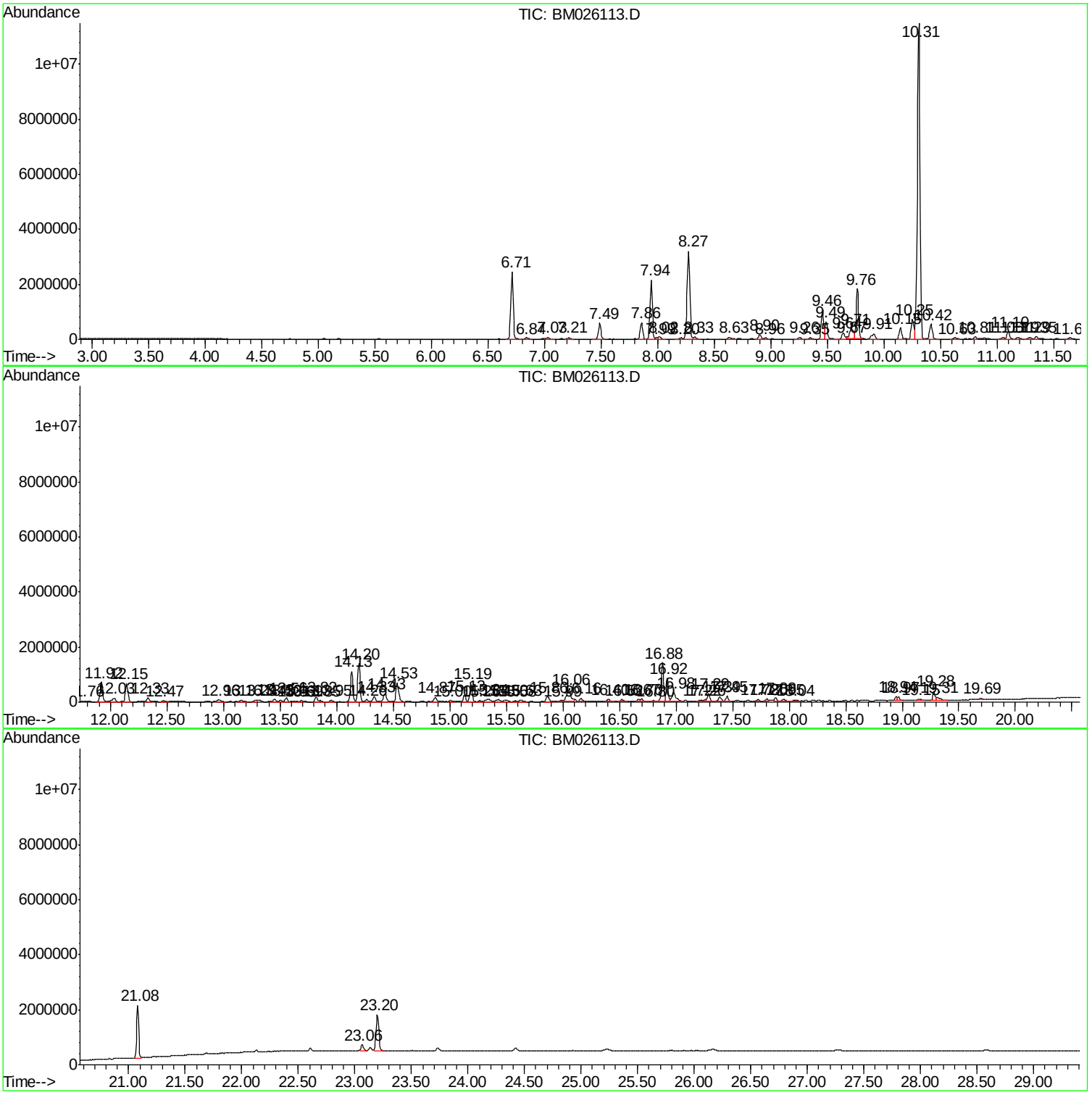
LSC Report - Integrated Chromatogram

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM052620\
Data File : BM026113.D
Acq On : 26 May 2020 14:26
Operator : MA/SJ
Sample : L2737-16DL 10X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9B13DL

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\BM052620\
Data File : BM026113.D
Acq On : 26 May 2020 14:26
Operator : MA/SJ
Sample : L2737-16DL 10X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9B13DL

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

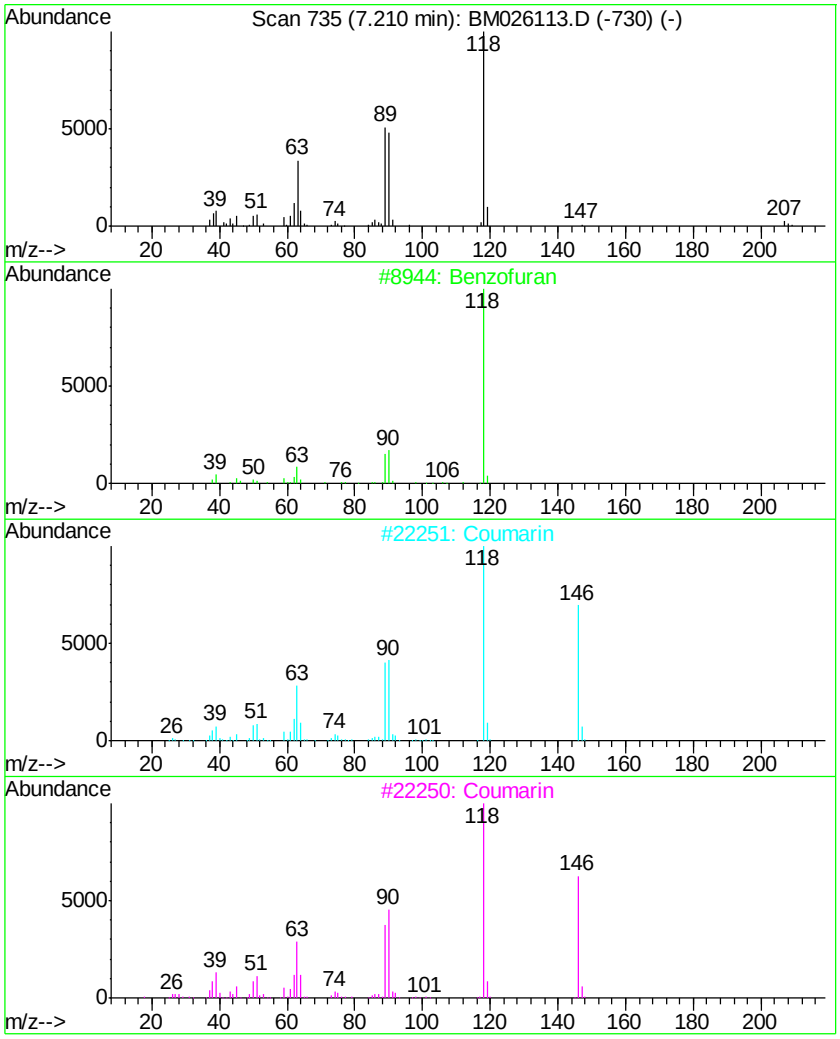
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

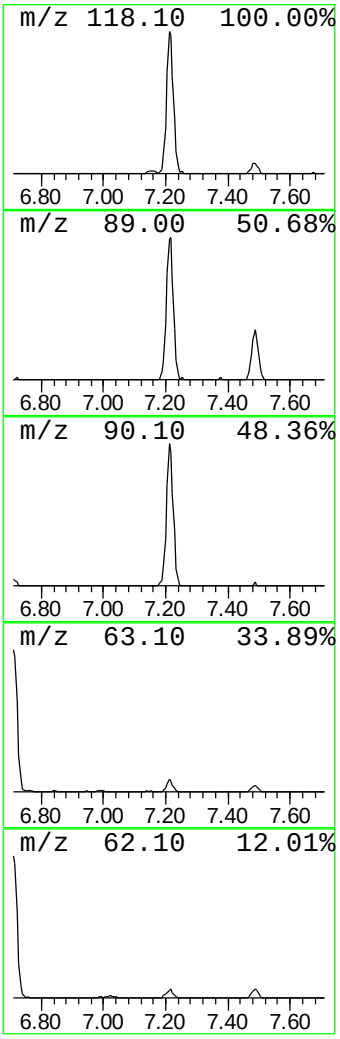
****

Peak Number 1 Benzofuran Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.21	2.06 ng/ul	92872	1,4-Dichlorobenzene-d4	7.49

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzofuran	118	C8H6O	000271-89-6	90
2	Coumarin	146	C9H6O2	000091-64-5	90
3	Coumarin	146	C9H6O2	000091-64-5	83
4	Benzofuran	118	C8H6O	000271-89-6	76
5	3-Hydroxyphenylacetylene	118	C8H6O	010401-11-3	72





Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM052620\
Data File : BM026113.D
Acq On : 26 May 2020 14:26
Operator : MA/SJ
Sample : L2737-16DL 10X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9B13DL

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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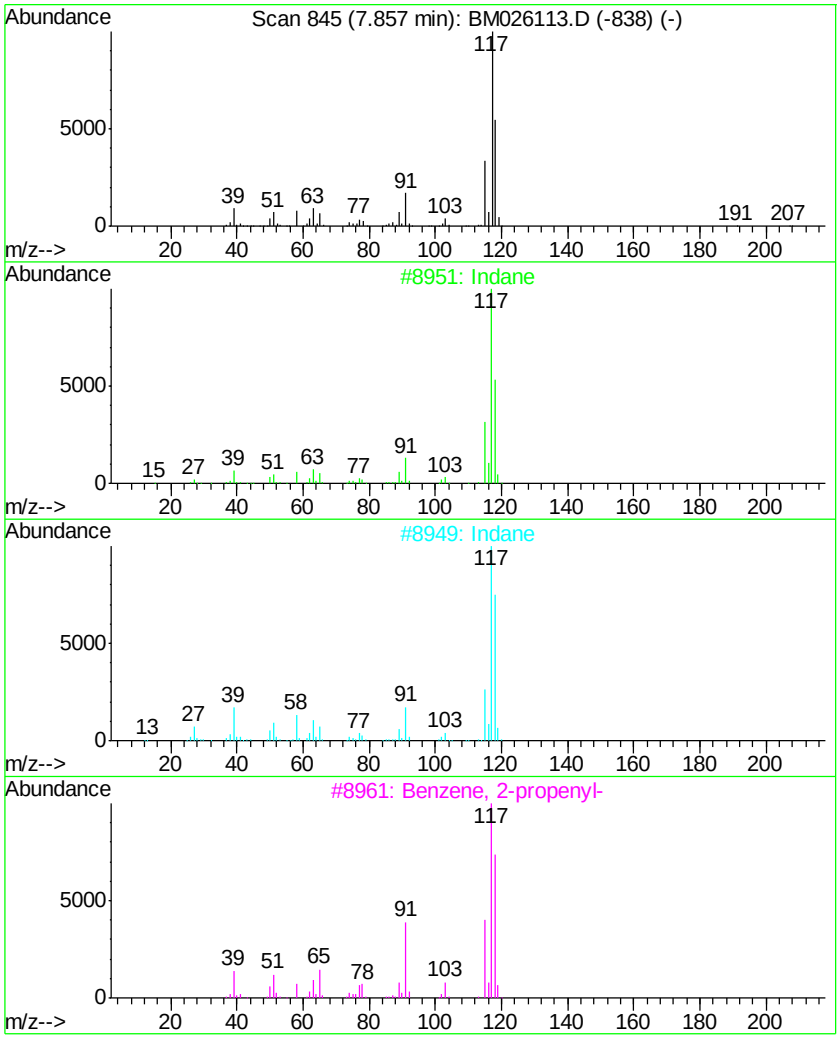
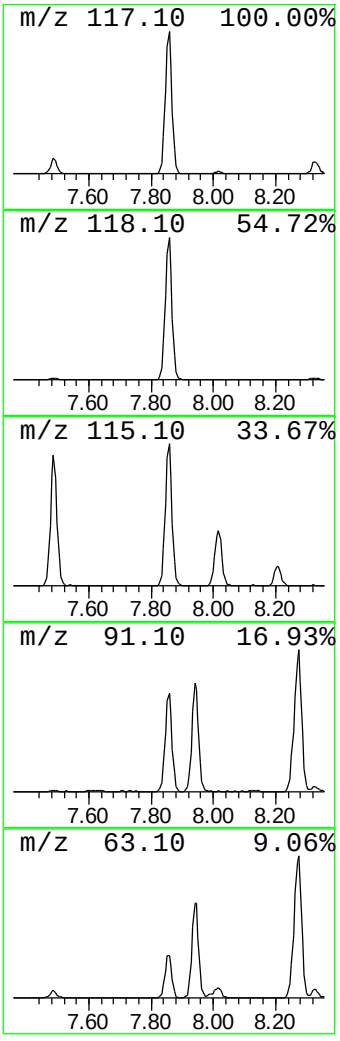
Peak Number 2 Indane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.86	20.49 ng/ul	925098	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	Benzene, 2-propenyl-	118	C9H10	000300-57-2	81
4	Indane	118	C9H10	000496-11-7	81
5	Tetracyclo[3.3.1.0(2,8).0(4,6)]-...	118	C9H10	1000191-13-7	80





Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM052620\
Data File : BM026113.D
Acq On : 26 May 2020 14:26
Operator : MA/SJ
Sample : L2737-16DL 10X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9B13DL

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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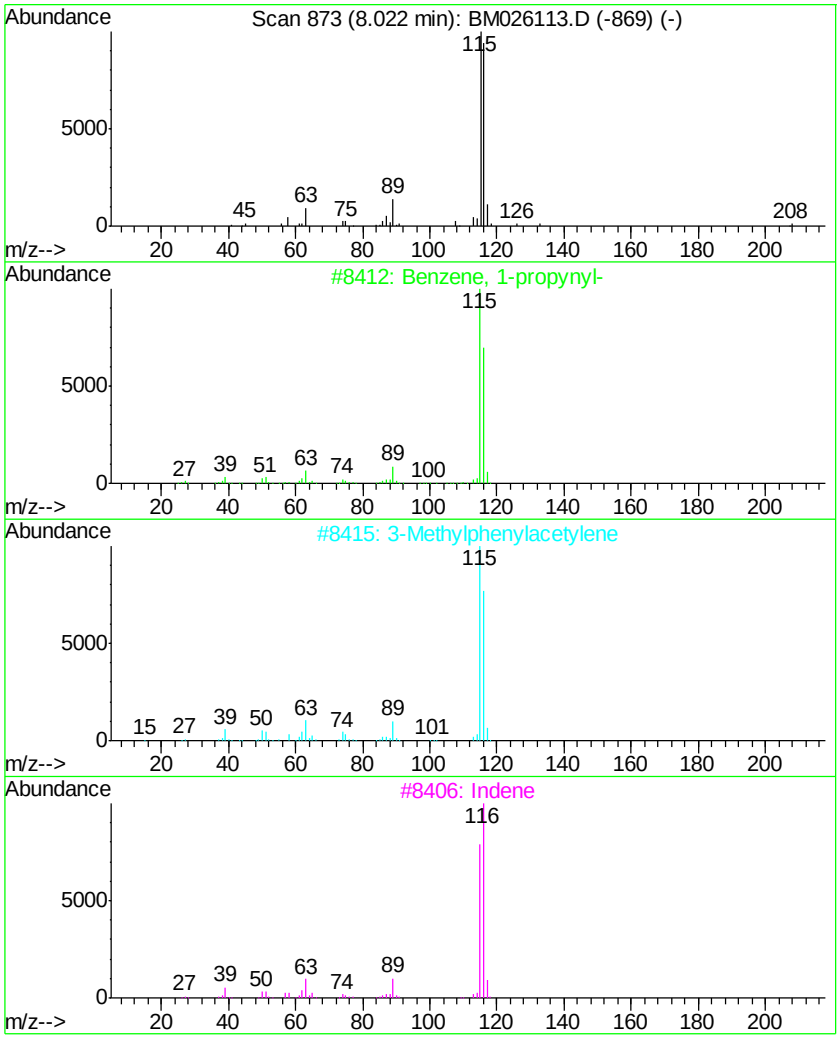
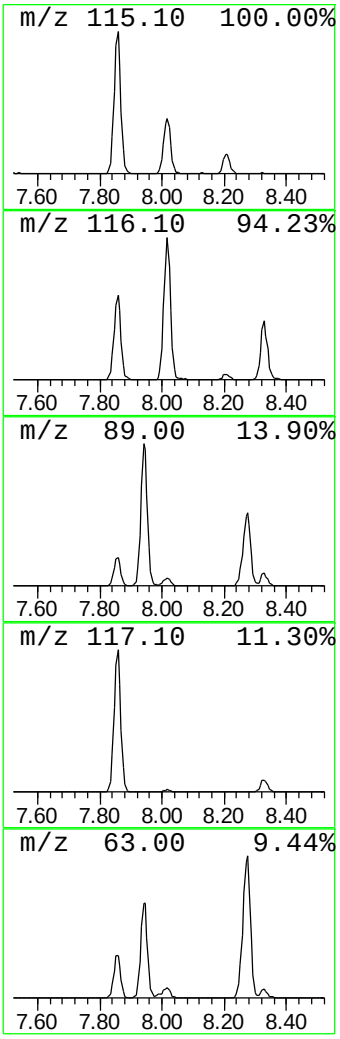
Peak Number 3 Benzene, 1-propynyl- Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.02	2.93 ng/ul	132443	1,4-Dichlorobenzene-d4	7.49

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-propynyl-	116	C9H8	000673-32-5	91
2	3-Methylphenylacetylene	116	C9H8	000766-82-5	91
3	Indene	116	C9H8	000095-13-6	90
4	Benzene, 1-ethynyl-4-methyl-	116	C9H8	000766-97-2	90
5	Indene	116	C9H8	000095-13-6	86





Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM052620\
Data File : BM026113.D
Acq On : 26 May 2020 14:26
Operator : MA/SJ
Sample : L2737-16DL 10X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9B13DL

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

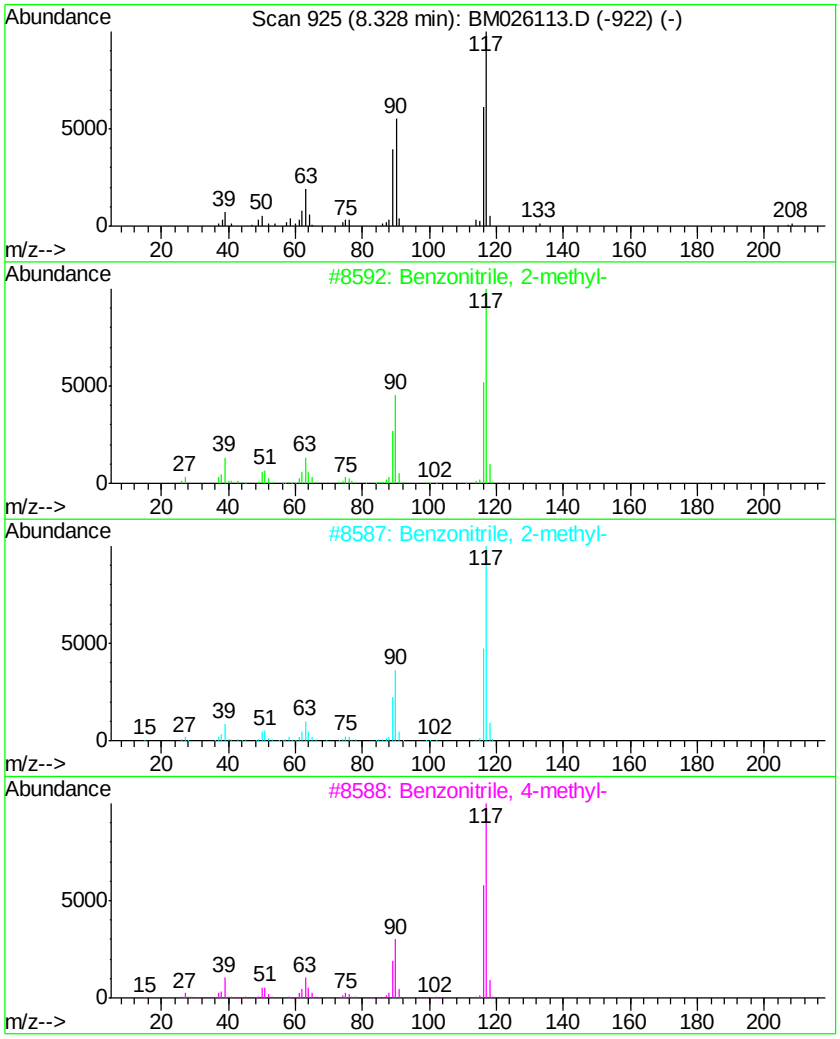
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

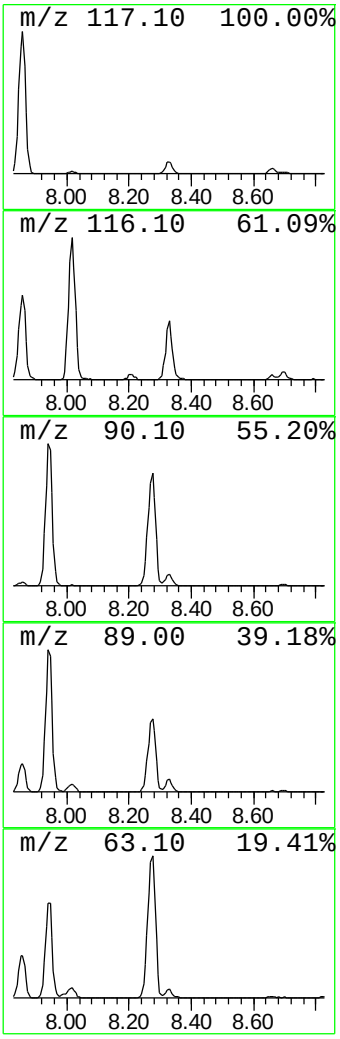
****

Peak Number 4 Benzonitrile, 2-methyl- Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.33	2.57 ng/ul	116055	1,4-Dichlorobenzene-d4	7.49

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzonitrile, 2-methyl-	117	C8H7N	000529-19-1	91
2	Benzonitrile, 2-methyl-	117	C8H7N	000529-19-1	90
3	Benzonitrile, 4-methyl-	117	C8H7N	000104-85-8	90
4	Benzonitrile, 3-methyl-	117	C8H7N	000620-22-4	87
5	Benzene, 1-isocyano-2-methyl-	117	C8H7N	010468-64-1	87





Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM052620\
Data File : BM026113.D
Acq On : 26 May 2020 14:26
Operator : MA/SJ
Sample : L2737-16DL 10X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9B13DL

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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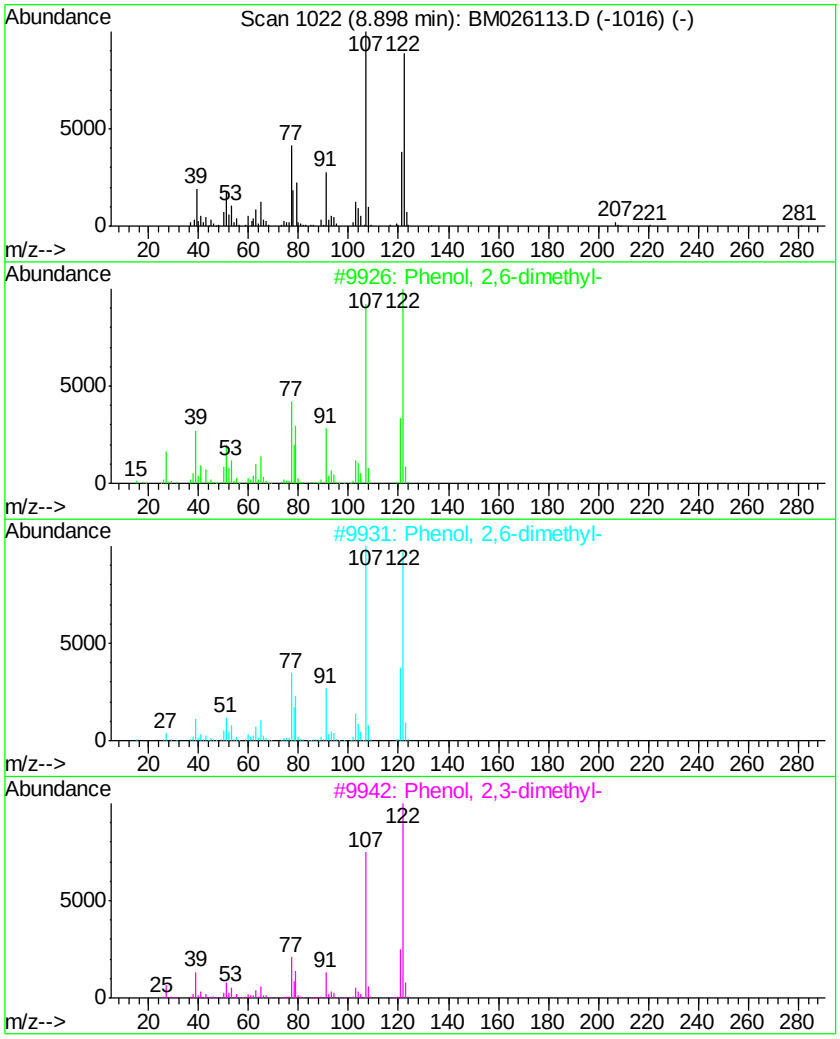
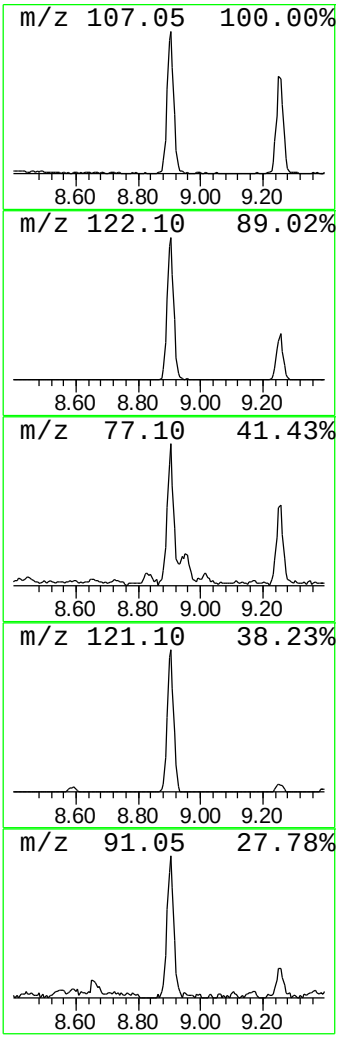
Peak Number 5 Phenol, 2,6-dimethyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.90	4.45 ng/ul	272233	Naphthalene-d8	10.25

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 2,6-dimethyl-	122	C8H10O	000576-26-1	98
2	Phenol, 2,6-dimethyl-	122	C8H10O	000576-26-1	97
3	Phenol, 2,3-dimethyl-	122	C8H10O	000526-75-0	95
4	Phenol, 2,5-dimethyl-	122	C8H10O	000095-87-4	95
5	Phenol, 2,5-dimethyl-	122	C8H10O	000095-87-4	94





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Data File : BM026113.D
Acq On : 26 May 2020 14:26
Operator : MA/SJ
Sample : L2737-16DL 10X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9B13DL

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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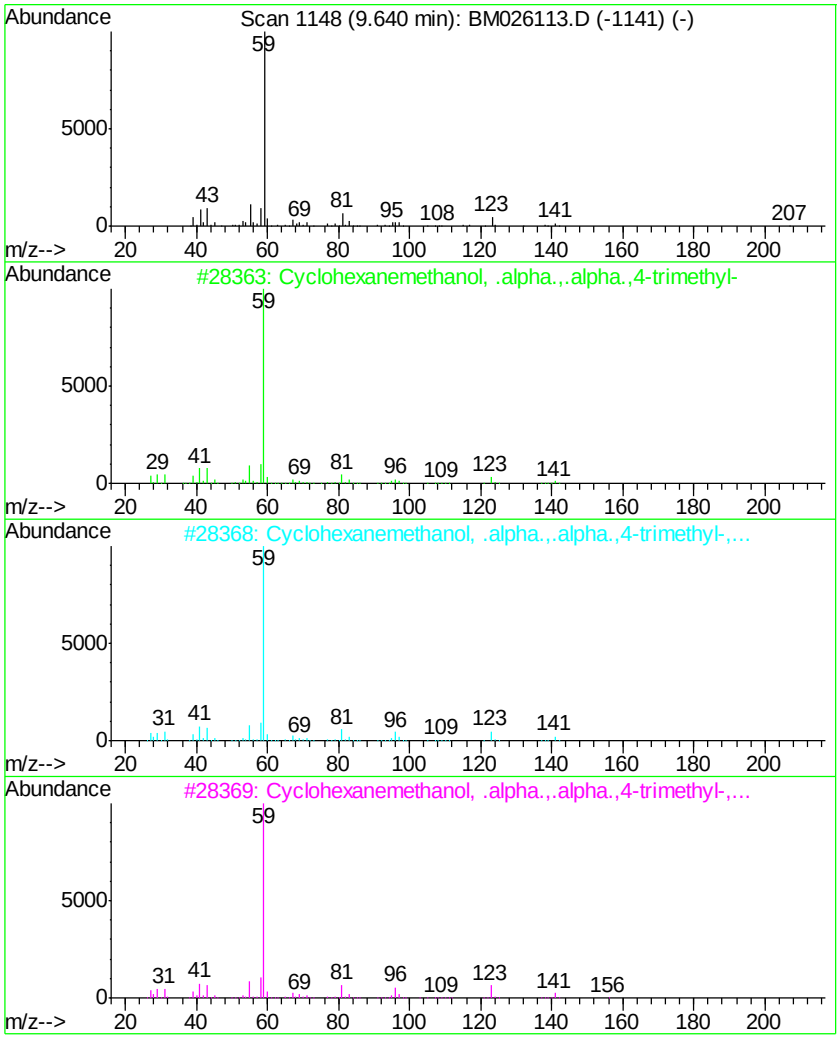
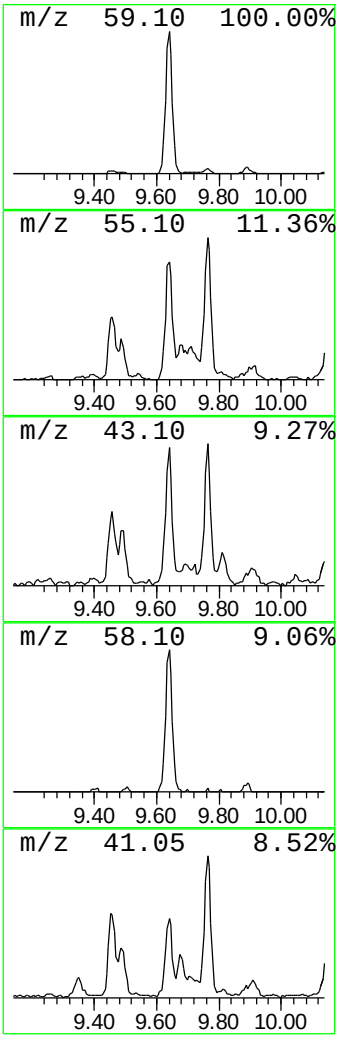
Peak Number 6 Cyclohexanemethanol, .alpha... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.64	6.41 ng/ul	392085	Naphthalene-d8	10.25

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexanemethanol, .alpha.,.al...	156	C10H20O	000498-81-7	83
2	Cyclohexanemethanol, .alpha.,.al...	156	C10H20O	007322-63-6	78
3	Cyclohexanemethanol, .alpha.,.al...	156	C10H20O	005114-00-1	78
4	Butane, 2-methoxy-3-methyl-	102	C6H14O	062016-49-3	50
5	Cyclohexanemethanol, .alpha.,.al...	156	C10H20O	000498-81-7	47





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Acq On : 26 May 2020 14:26
Operator : MA/SJ
Sample : L2737-16DL 10X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9B13DL

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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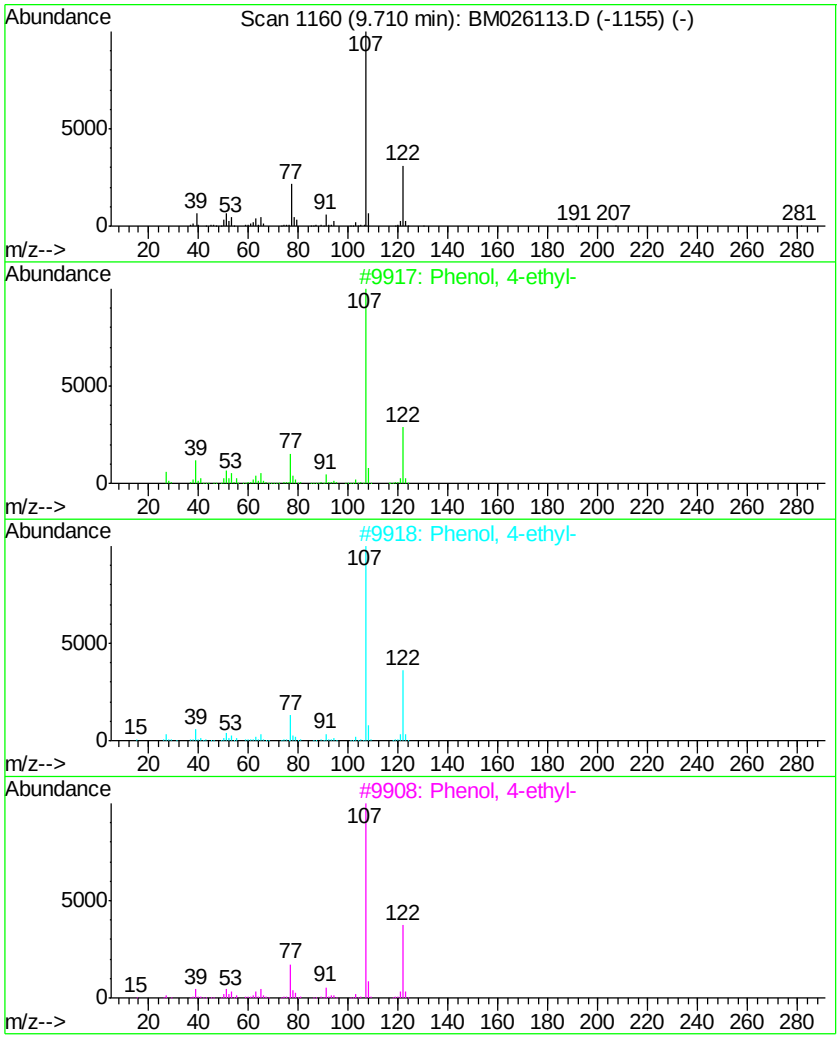
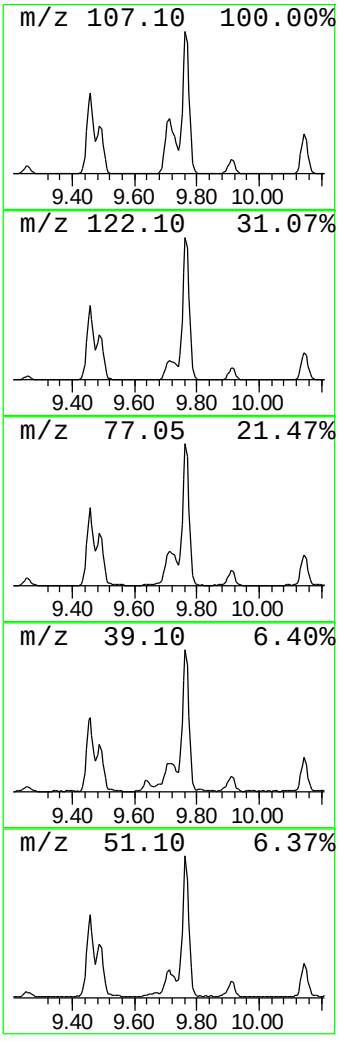
Peak Number 7 Phenol, 4-ethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.71	14.68 ng/ul	897270	Naphthalene-d8	10.25

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 4-ethyl-	122	C8H10O	000123-07-9	93
2	Phenol, 4-ethyl-	122	C8H10O	000123-07-9	91
3	Phenol, 4-ethyl-	122	C8H10O	000123-07-9	91
4	Phenol, 2-ethyl-	122	C8H10O	000090-00-6	91
5	Phenol, 2-ethyl-	122	C8H10O	000090-00-6	91





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM052620\
Data File : BM026113.D
Acq On : 26 May 2020 14:26
Operator : MA/SJ
Sample : L2737-16DL 10X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9B13DL

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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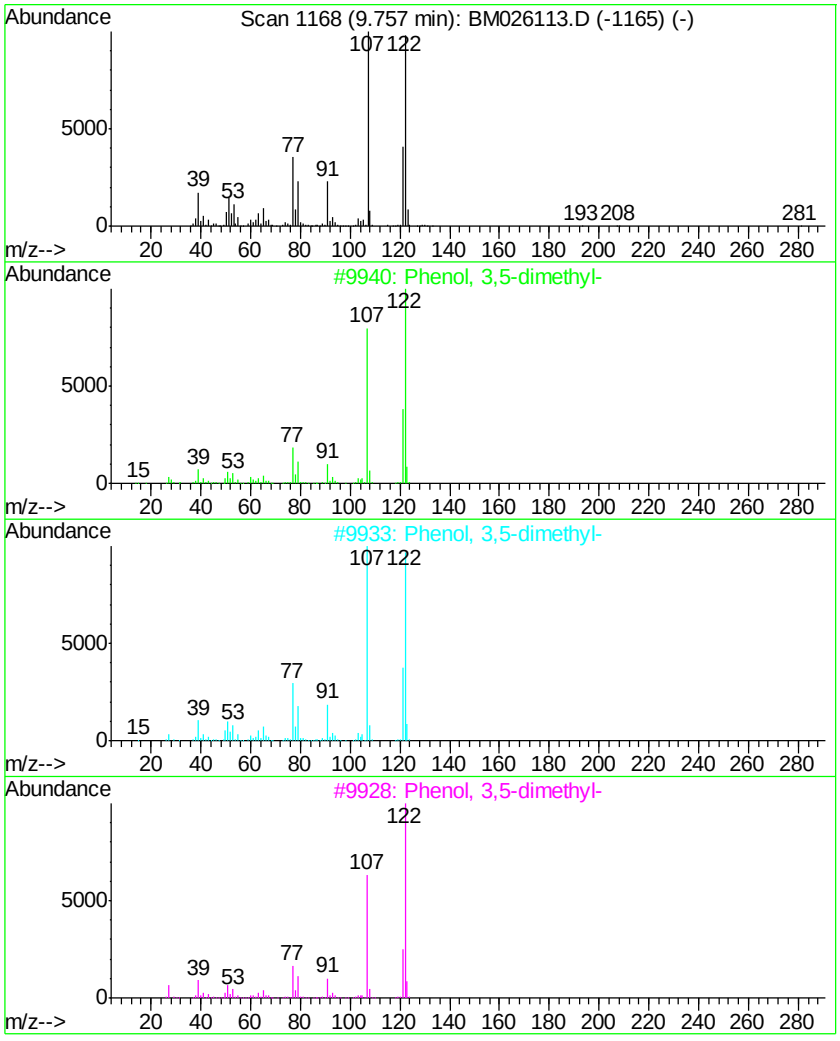
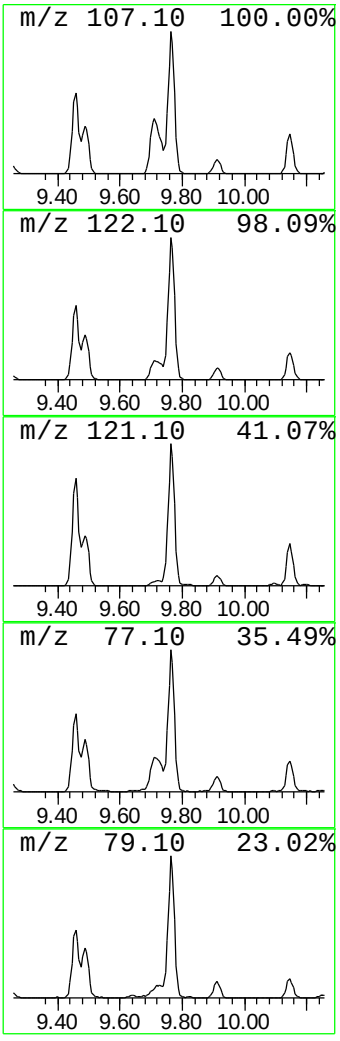
Peak Number 8 Phenol, 3,5-dimethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.76	43.24 ng/ul	2643140	Naphthalene-d8	10.25

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 3,5-dimethyl-	122	C8H10O	000108-68-9	95
2	Phenol, 3,5-dimethyl-	122	C8H10O	000108-68-9	94
3	Phenol, 3,5-dimethyl-	122	C8H10O	000108-68-9	93
4	Phenol, 2,5-dimethyl-	122	C8H10O	000095-87-4	93
5	Phenol, 3,4-dimethyl-	122	C8H10O	000095-65-8	90





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Data File : BM026113.D
Acq On : 26 May 2020 14:26
Operator : MA/SJ
Sample : L2737-16DL 10X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9B13DL

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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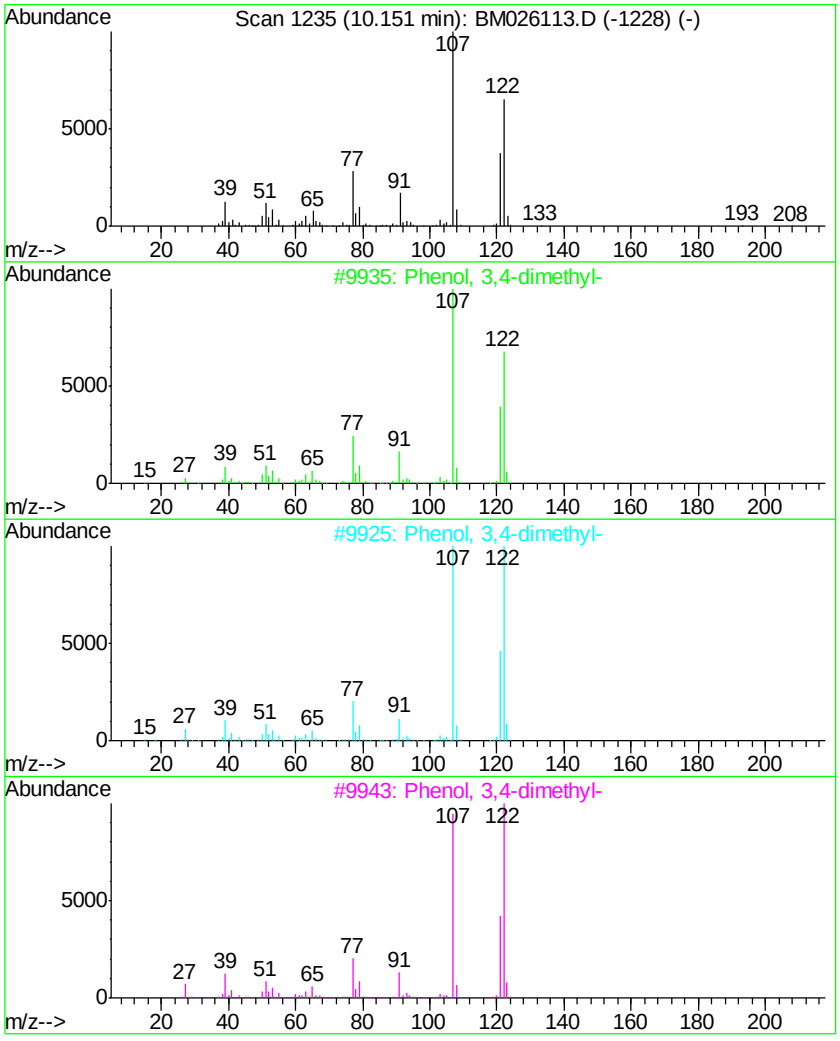
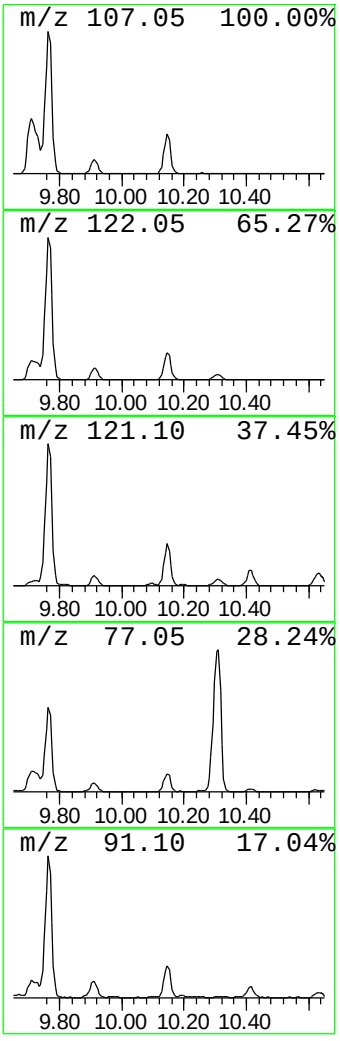
Peak Number 9 Phenol, 3,4-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.15	10.58 ng/ul	646503	Naphthalene-d8	10.25

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 3,4-dimethyl-	122	C8H10O	000095-65-8	96
2	Phenol, 3,4-dimethyl-	122	C8H10O	000095-65-8	95
3	Phenol, 3,4-dimethyl-	122	C8H10O	000095-65-8	95
4	Phenol, 3-ethyl-	122	C8H10O	000620-17-7	91
5	Phenol, 2-ethyl-	122	C8H10O	000090-00-6	91





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Acq On : 26 May 2020 14:26
Operator : MA/SJ
Sample : L2737-16DL 10X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9B13DL

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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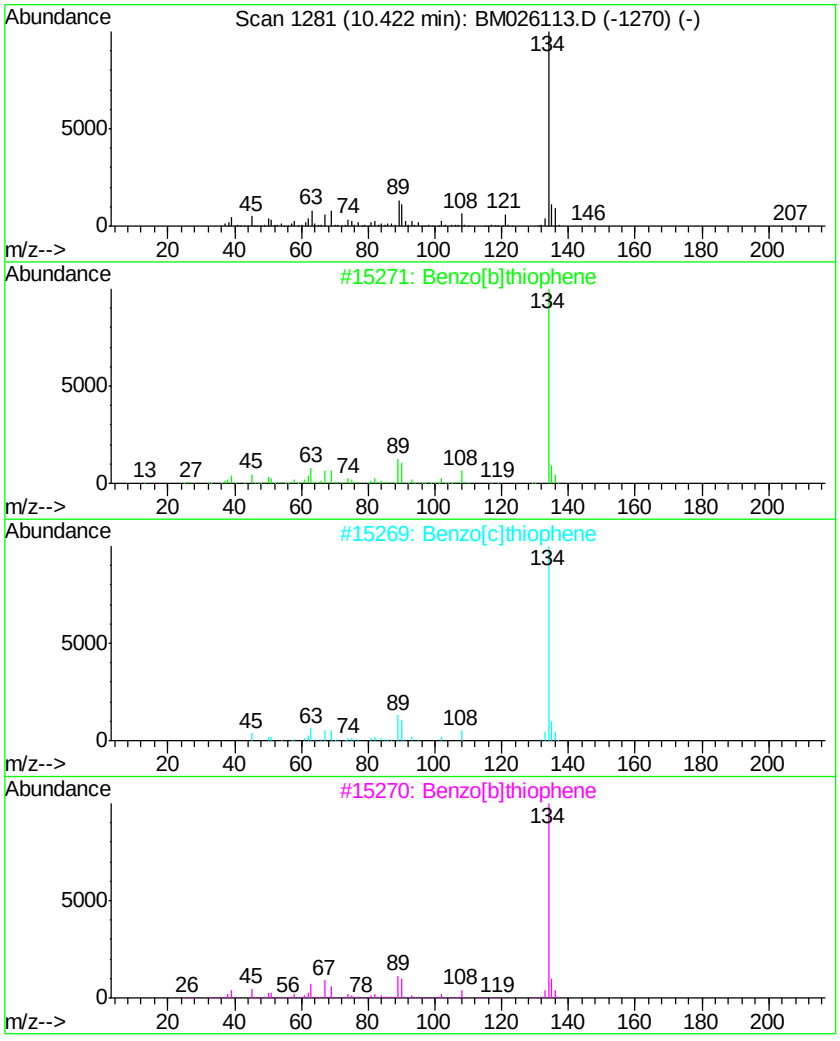
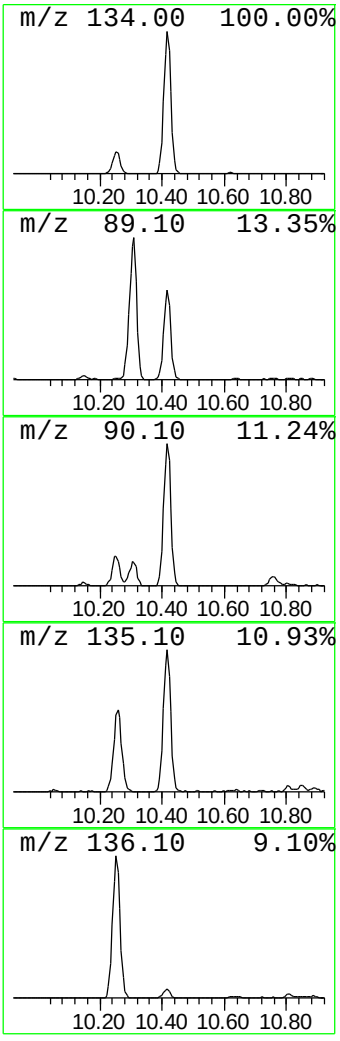
Peak Number 10 Benzo[b]thiophene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.42	15.15 ng/ul	926060	Naphthalene-d8	10.25

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzo[b]thiophene	134	C8H6S	000095-15-8	94
2	Benzo[c]thiophene	134	C8H6S	000270-82-6	93
3	Benzo[b]thiophene	134	C8H6S	000095-15-8	93
4	Cyclopenta[c]thiapyran	134	C8H6S	000270-63-3	90
5	Benzo[b]thiophene	134	C8H6S	000095-15-8	87





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Operator : MA/SJ
Sample : L2737-16DL 10X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9B13DL

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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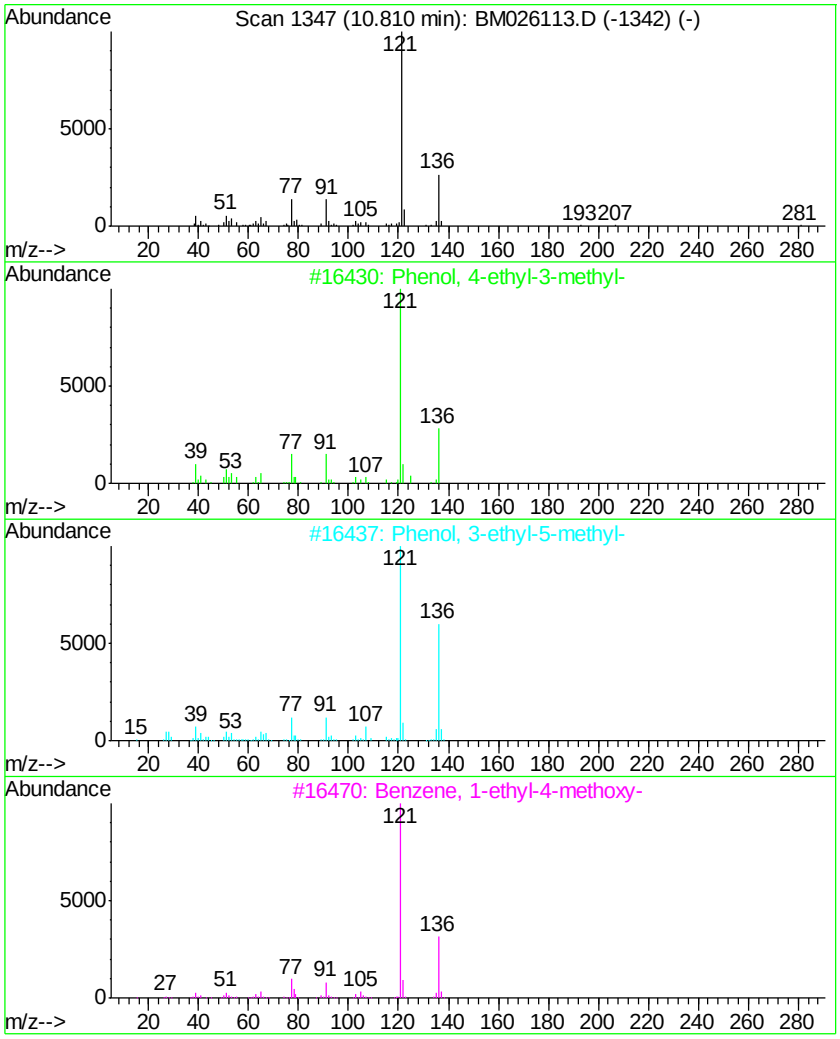
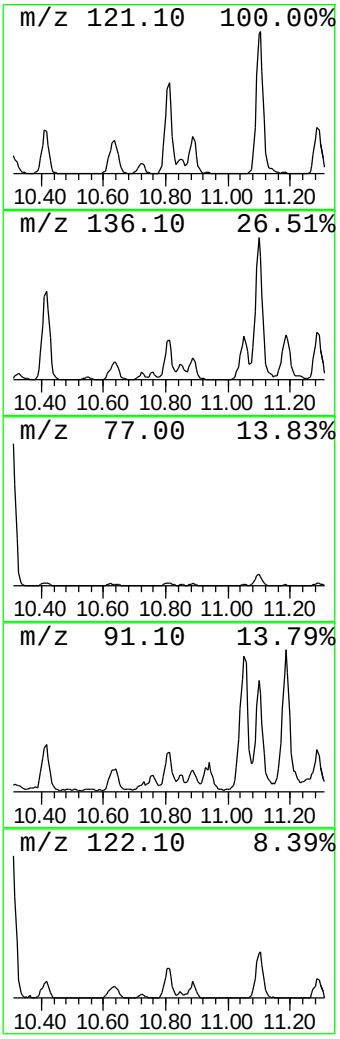
Peak Number 11 Phenol, 4-ethyl-3-methyl- Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.81	2.03 ng/ul	124125	Naphthalene-d8	10.25

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 4-ethyl-3-methyl-	136	C9H12O	001123-94-0	91
2	Phenol, 3-ethyl-5-methyl-	136	C9H12O	000698-71-5	91
3	Benzene, 1-ethyl-4-methoxy-	136	C9H12O	001515-95-3	91
4	Phenol, 2-(1-methylethyl)-	136	C9H12O	000088-69-7	91
5	Phenol, 3-(1-methylethyl)-	136	C9H12O	000618-45-1	91





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Acq On : 26 May 2020 14:26
Operator : MA/SJ
Sample : L2737-16DL 10X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9B13DL

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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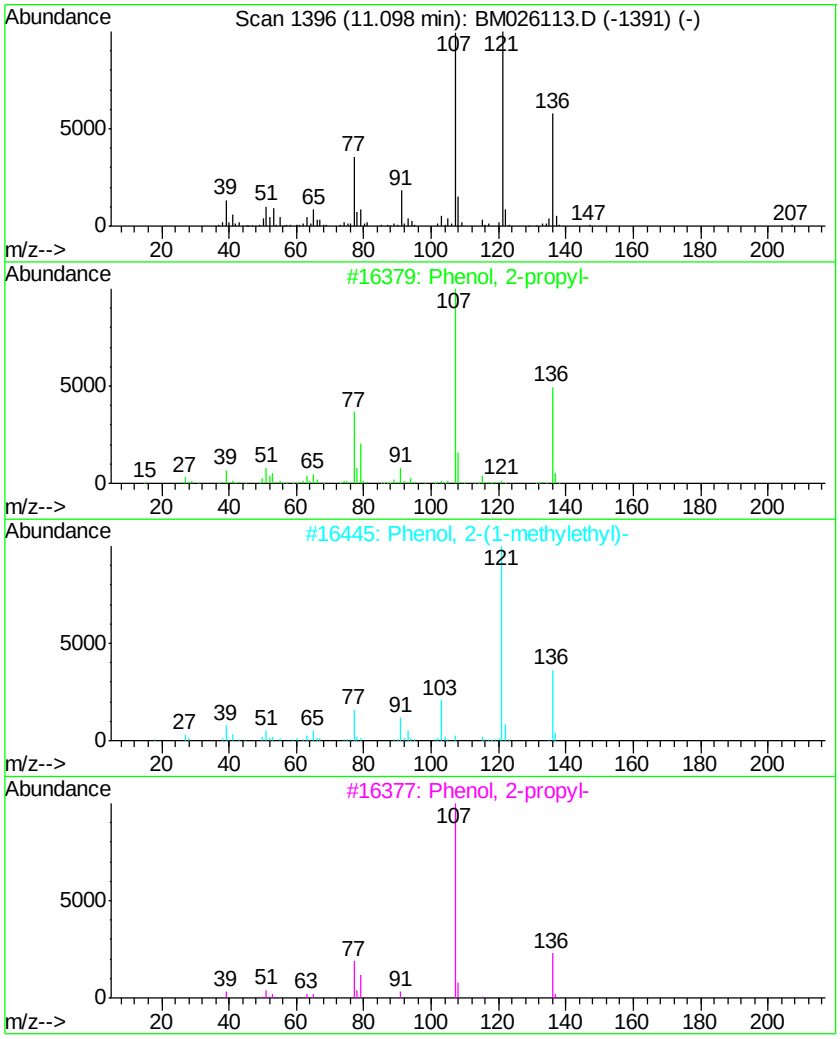
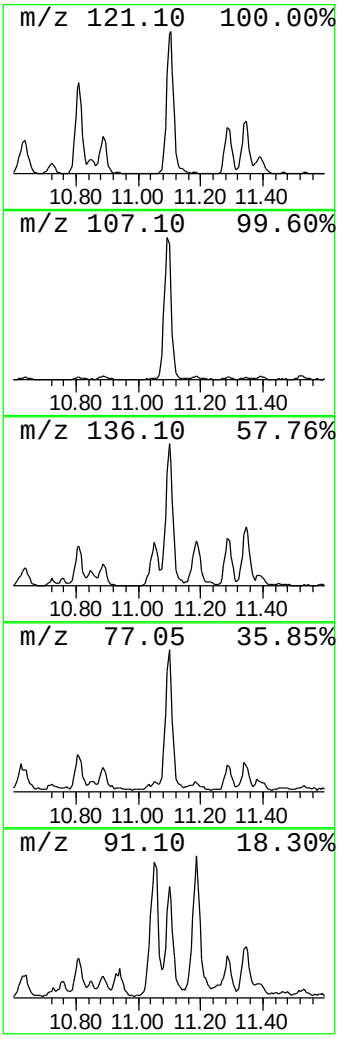
Peak Number 12 Phenol, 2-propyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.10	7.25 ng/ul	442923	Naphthalene-d8	10.25

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 2-propyl-	136	C9H12O	000644-35-9	53
2	Phenol, 2-(1-methylethyl)-	136	C9H12O	000088-69-7	53
3	Phenol, 2-propyl-	136	C9H12O	000644-35-9	53
4	Phenol, 3-ethyl-5-methyl-	136	C9H12O	000698-71-5	53
5	Benzene, 1-ethyl-4-methoxy-	136	C9H12O	001515-95-3	52





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Operator : MA/SJ
Sample : L2737-16DL 10X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9B13DL

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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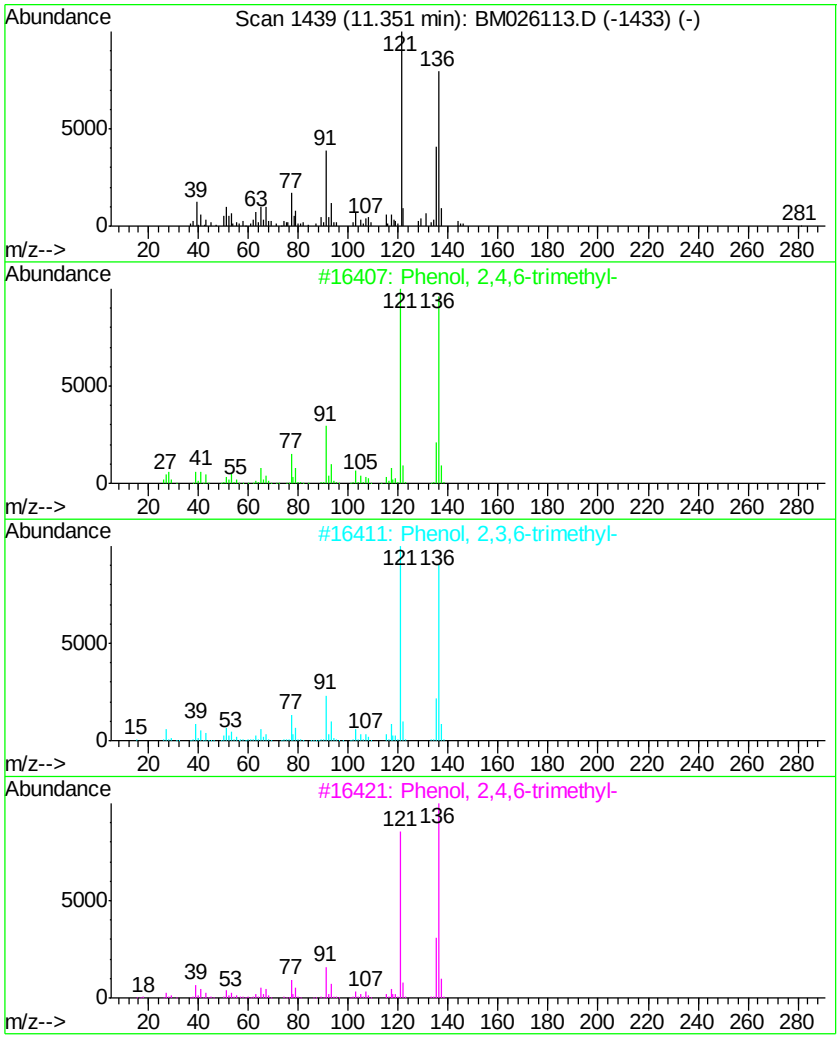
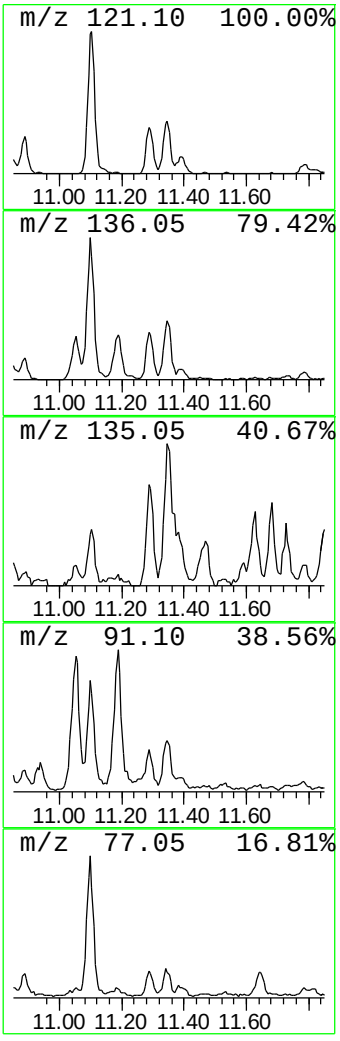
Peak Number 13 Phenol, 2,4,6-trimethyl- Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.35	2.23 ng/ul	136496	Naphthalene-d8	10.25

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 2,4,6-trimethyl-	136	C9H12O	000527-60-6	95
2	Phenol, 2,3,6-trimethyl-	136	C9H12O	002416-94-6	94
3	Phenol, 2,4,6-trimethyl-	136	C9H12O	000527-60-6	93
4	Phenol, 2,3,5-trimethyl-	136	C9H12O	000697-82-5	93
5	Phenol, 2,3,5-trimethyl-	136	C9H12O	000697-82-5	93





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Sample : L2737-16DL 10X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9B13DL

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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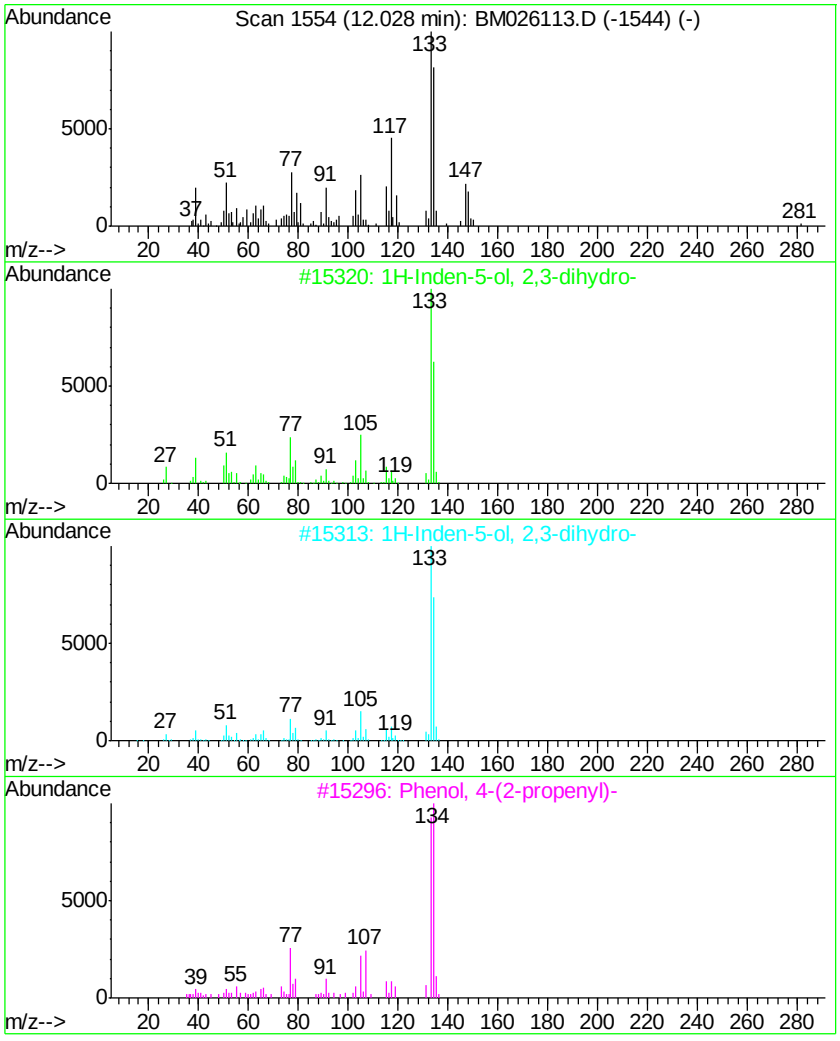
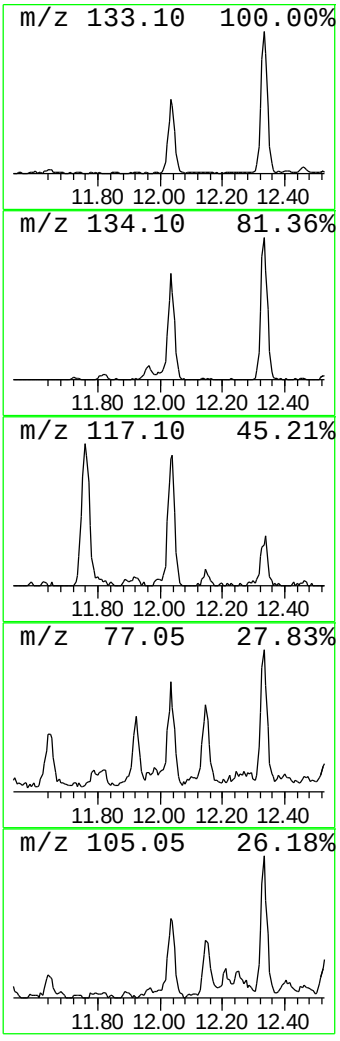
Peak Number 14 1H-Inden-5-ol, 2,3-dihydro- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.03	5.16 ng/ul	315658	Naphthalene-d8	10.25

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Inden-5-ol, 2,3-dihydro-	134	C9H10O	001470-94-6	86
2	1H-Inden-5-ol, 2,3-dihydro-	134	C9H10O	001470-94-6	70
3	Phenol, 4-(2-propenyl)-	134	C9H10O	000501-92-8	62
4	Benzaldehyde, 3,4-dimethyl-	134	C9H10O	005973-71-7	60
5	Phenol, 4-(2-propenyl)-	134	C9H10O	000501-92-8	60





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Operator : MA/SJ
Sample : L2737-16DL 10X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9B13DL

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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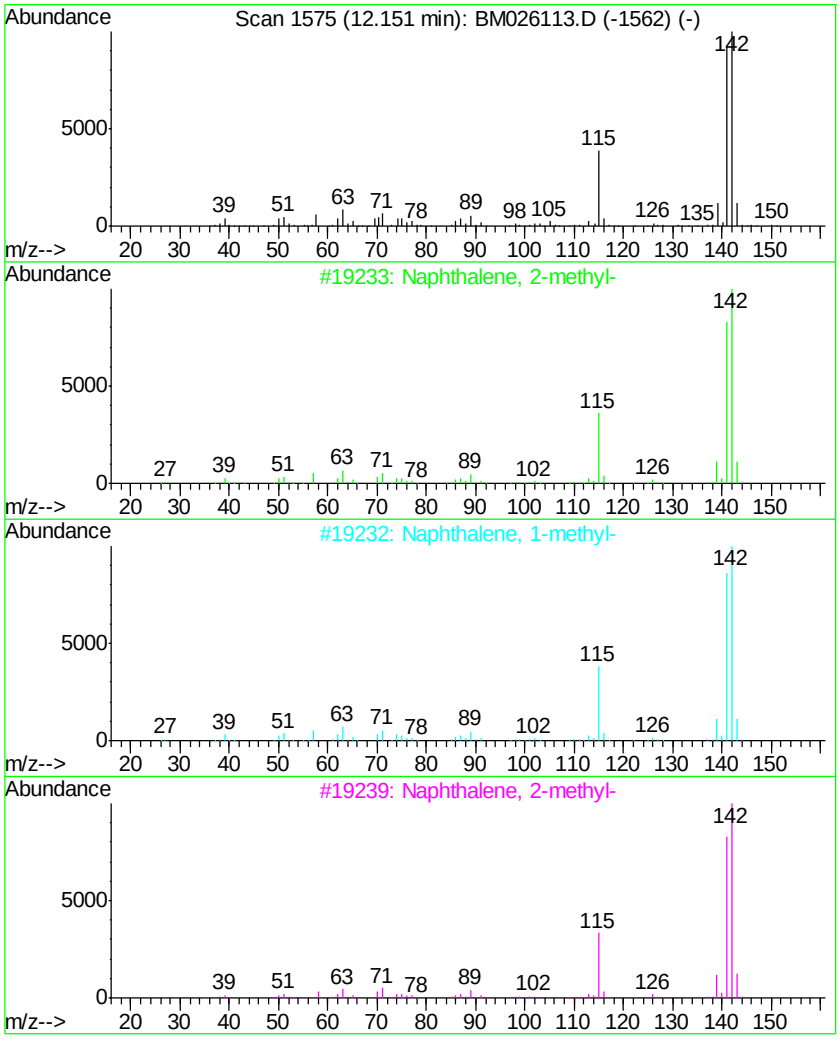
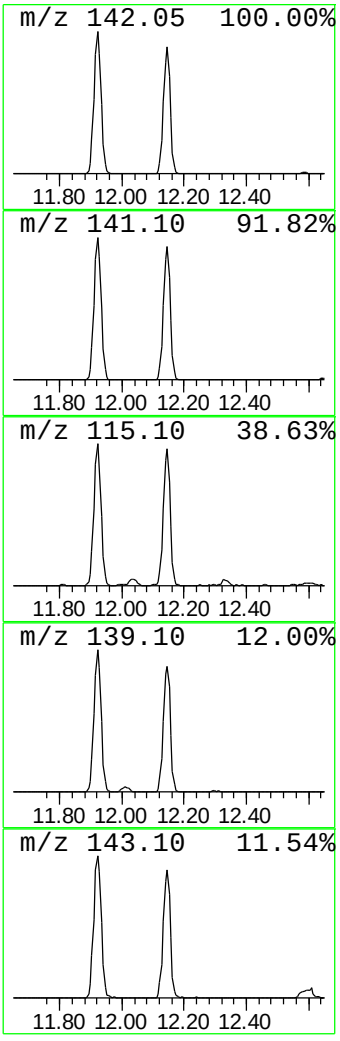
Peak Number 15 Naphthalene, 1-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.15	17.55 ng/ul	1072670	Naphthalene-d8	10.25

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
2	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
3	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
4	1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	94
5	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	93





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Operator : MA/SJ
Sample : L2737-16DL 10X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9B13DL

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 16 Benzaldehyde, 3,4-dimethyl- Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.33	2.75 ng/ul	220943	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	95
2	1H-Inden-5-ol, 2,3-dihydro-	134	C9H10O	001470-94-6	87
3	Benzaldehyde, 3,4-dimethyl-	134	C9H10O	005973-71-7	87
4	Benzaldehyde, 2,4-dimethyl-	134	C9H10O	015764-16-6	74
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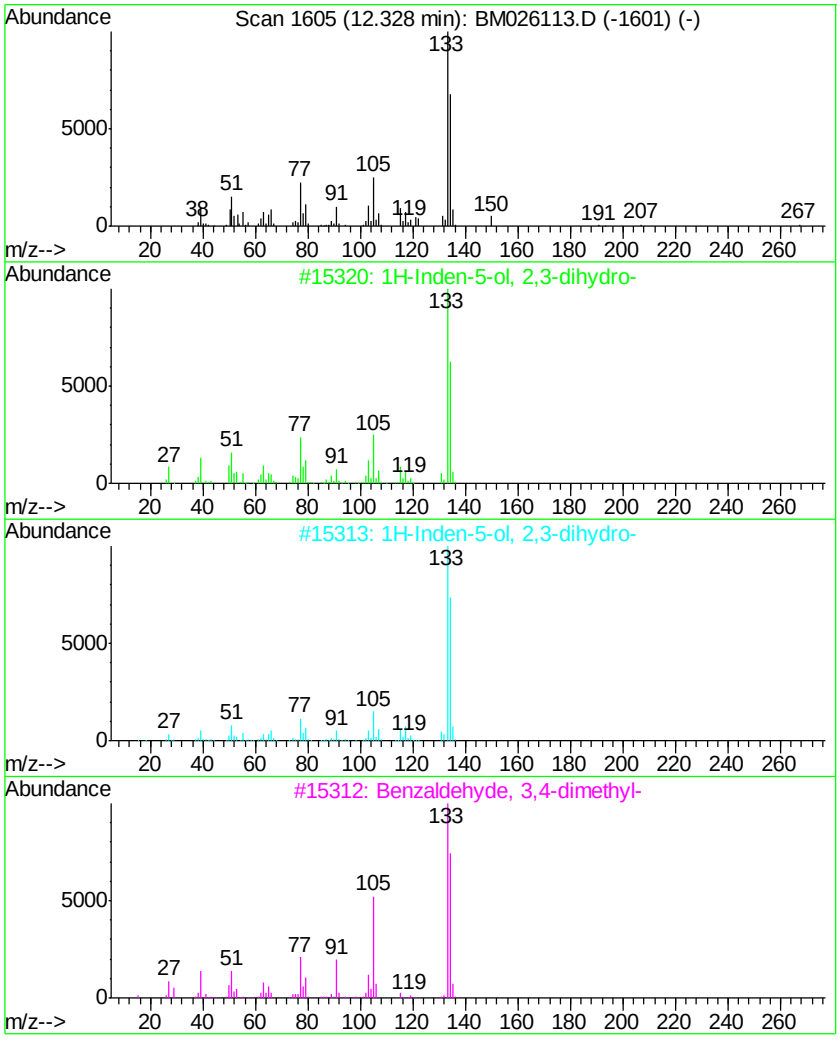

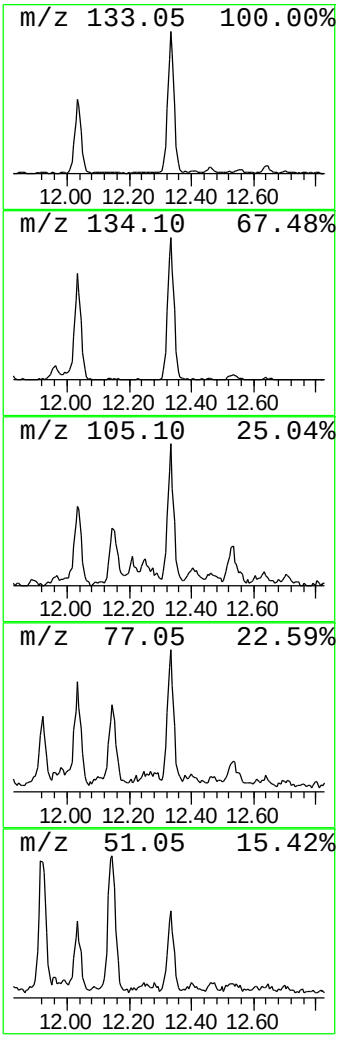

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Acq On : 26 May 2020 14:26
Operator : MA/SJ
Sample : L2737-16DL 10X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9B13DL

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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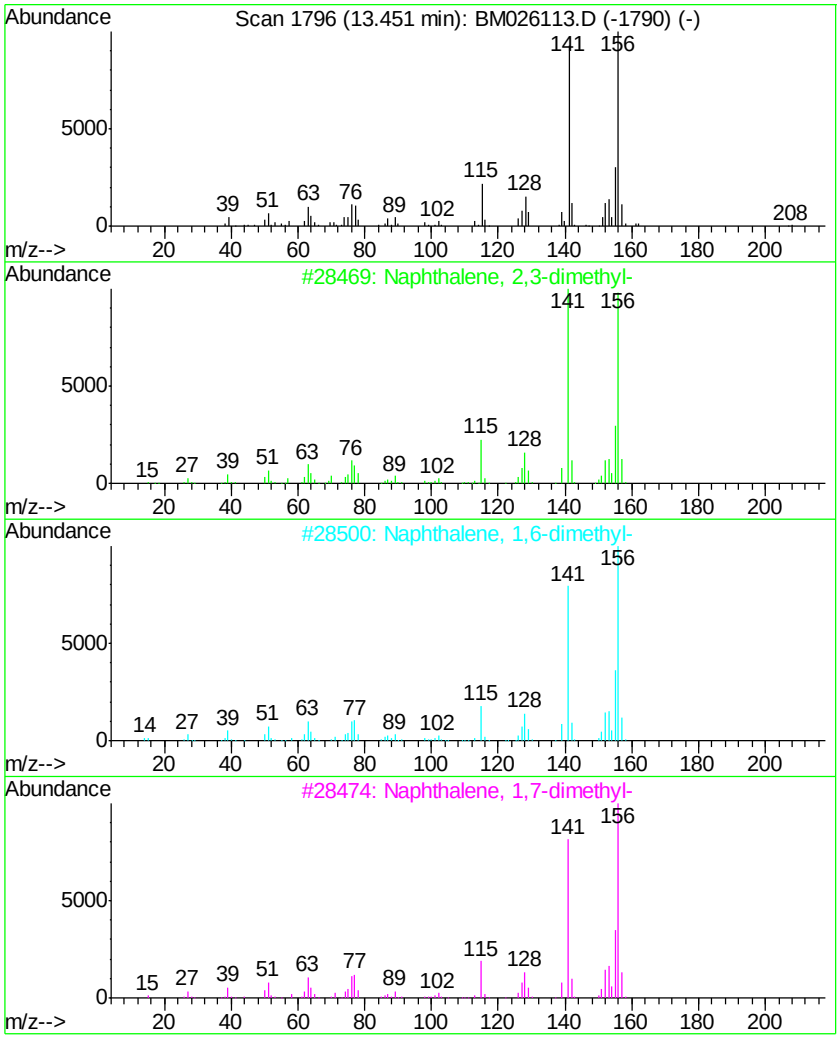
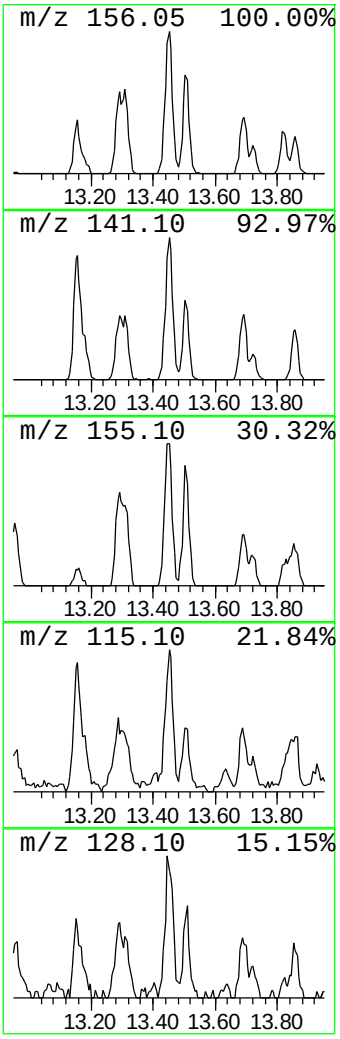
Peak Number 17 Naphthalene, 2,3-dimethyl- Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.45	2.37 ng/ul	190428	Acenaphthene-d10	14.13

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, 1,3-dimethyl-	156	C12H12	000575-41-7	97





Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM052620\
Data File : BM026113.D
Acq On : 26 May 2020 14:26
Operator : MA/SJ
Sample : L2737-16DL 10X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9B13DL

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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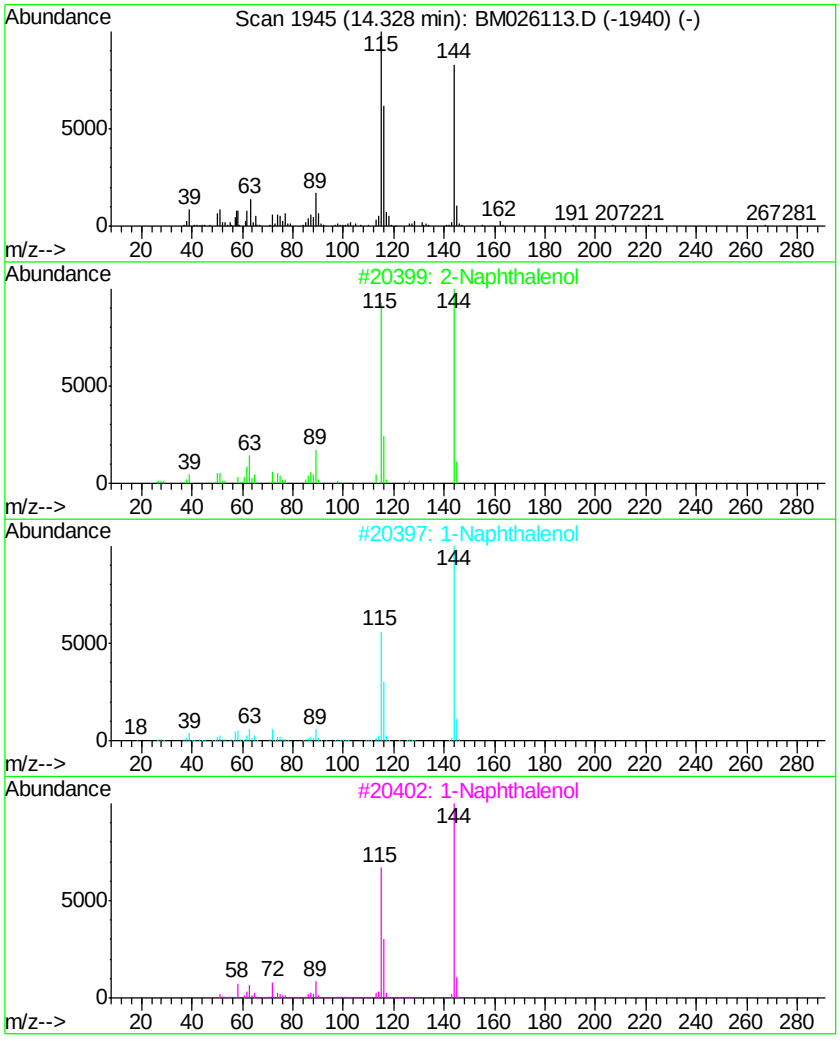
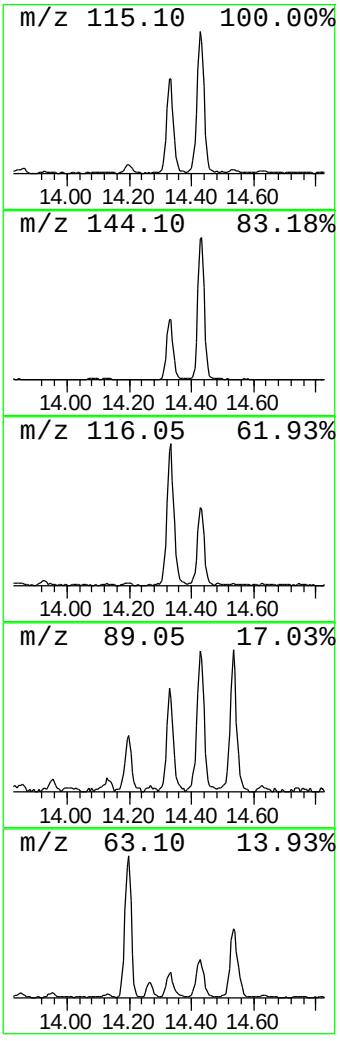
Peak Number 18 2-Naphthalenol Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.33	4.71 ng/ul	379108	Acenaphthene-d10	14.13

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Naphthalenol	144	C10H8O	000135-19-3	94
2	1-Naphthalenol	144	C10H8O	000090-15-3	93
3	1-Naphthalenol	144	C10H8O	000090-15-3	93
4	Benzene, 1-propynyl-	116	C9H8	000673-32-5	90
5	Benzene, 1-propynyl-	116	C9H8	000673-32-5	89





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Data File : BM026113.D
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Operator : MA/SJ
Sample : L2737-16DL 10X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9B13DL

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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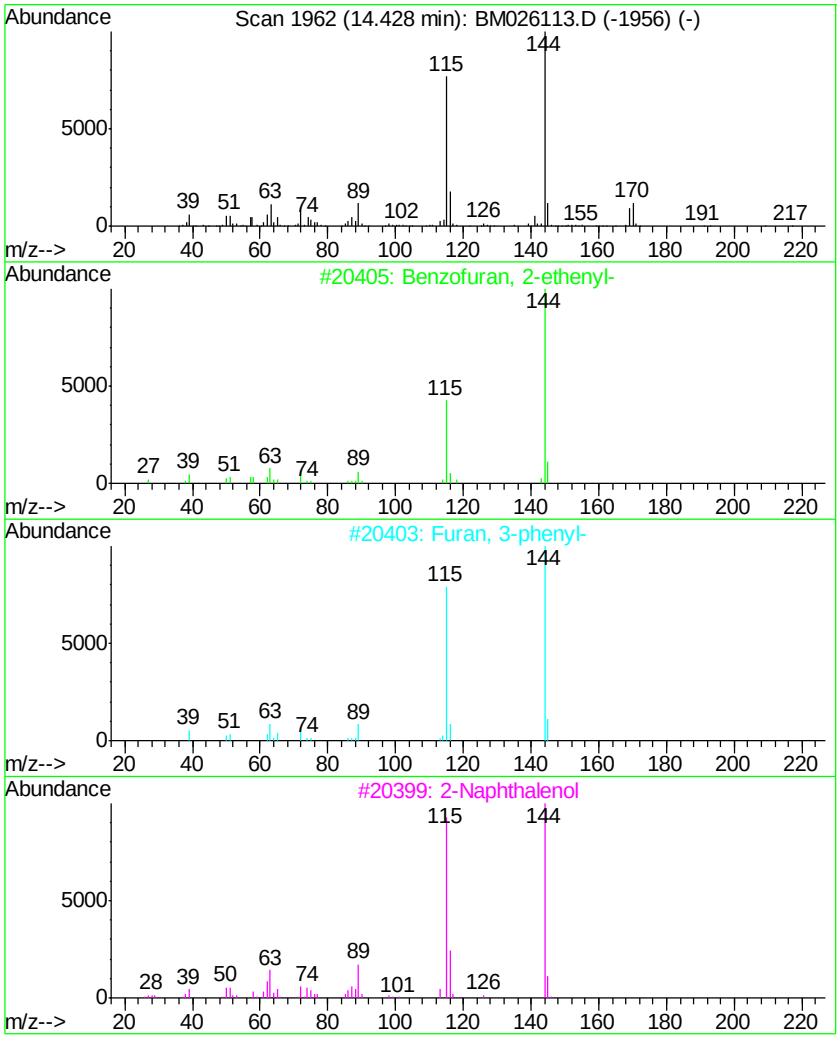
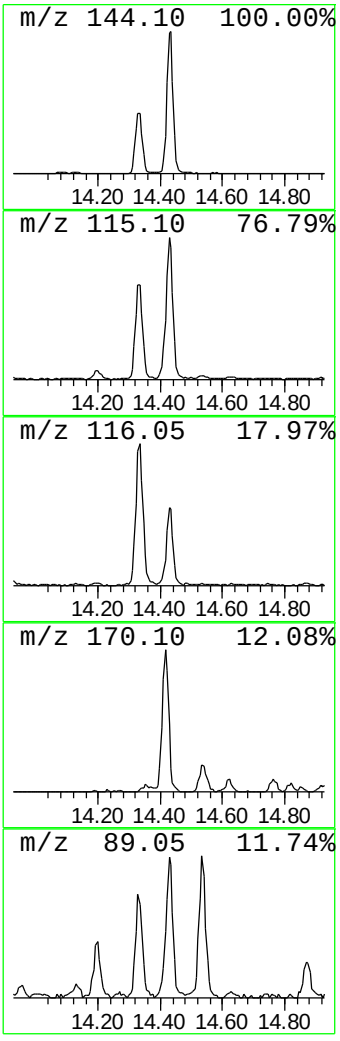
Peak Number 19 Benzofuran, 2-ethenyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.43	7.12 ng/ul	572610	Acenaphthene-d10	14.13

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzofuran, 2-ethenyl-	144	C10H8O	007522-79-4	87
2	Furan, 3-phenyl-	144	C10H8O	013679-41-9	87
3	2-Naphthalenol	144	C10H8O	000135-19-3	87
4	1-Naphthalenol	144	C10H8O	000090-15-3	83
5	1-Naphthalenol	144	C10H8O	000090-15-3	81





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Data File : BM026113.D
Acq On : 26 May 2020 14:26
Operator : MA/SJ
Sample : L2737-16DL 10X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9B13DL

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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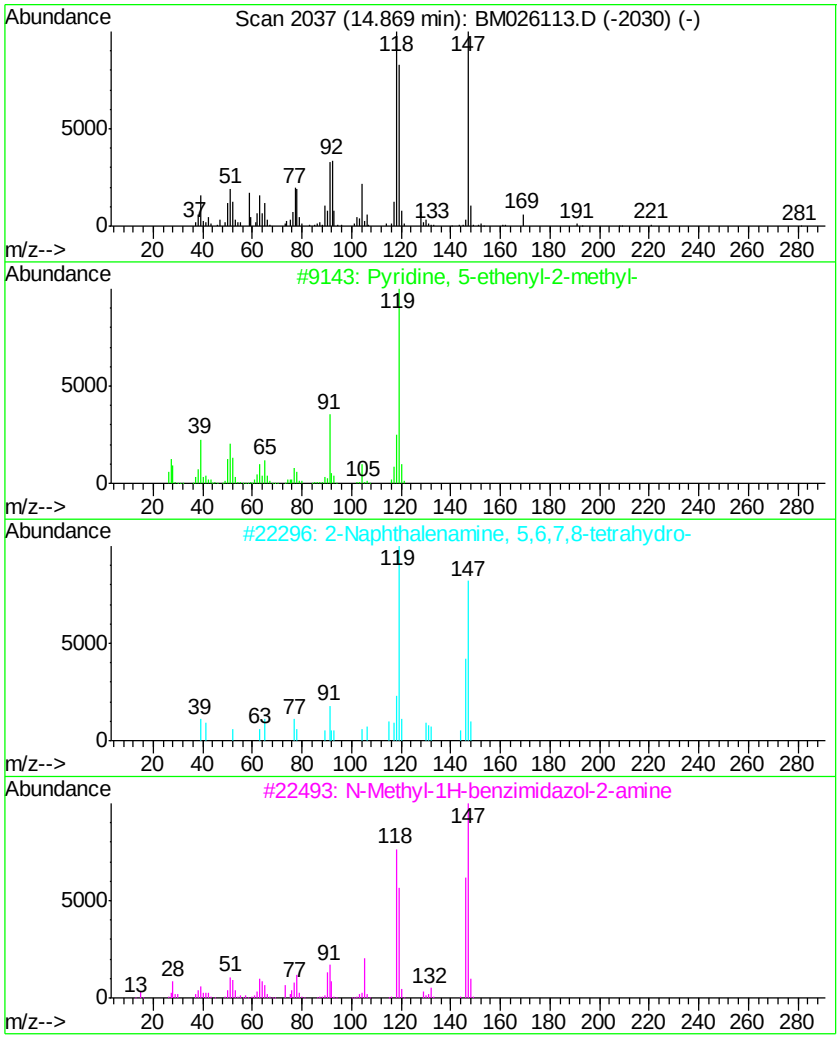
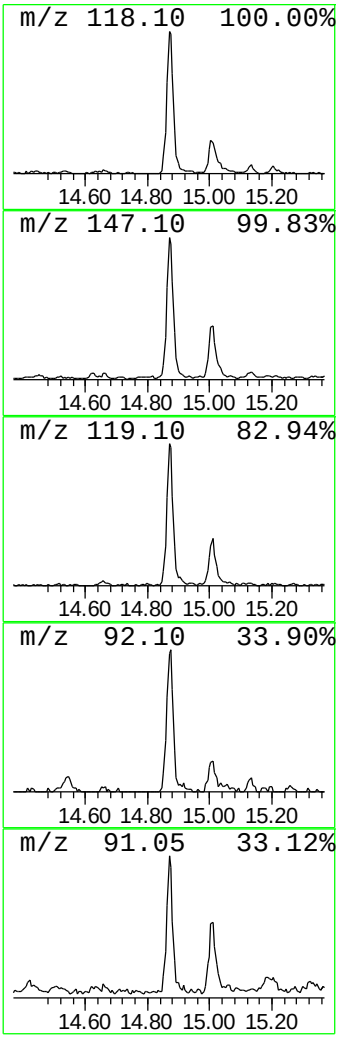
Peak Number 20 Pyridine, 5-ethenyl-2-methyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.87	3.09 ng/ul	248510	Acenaphthene-d10	14.13

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pyridine, 5-ethenyl-2-methyl-	119	C8H9N	000140-76-1	70
2	2-Naphthalenamine, 5,6,7,8-tetra...	147	C10H13N	002217-43-8	64
3	N-Methyl-1H-benzimidazol-2-amine	147	C8H9N3	1000332-71-6	64
4	Methanamine, N-(phenylmethylene)-	119	C8H9N	000622-29-7	58
5	7-Benzofuranamine, 2-methyl-	147	C9H9NO	1000327-46-4	58





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Acq On : 26 May 2020 14:26
Operator : MA/SJ
Sample : L2737-16DL 10X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9B13DL

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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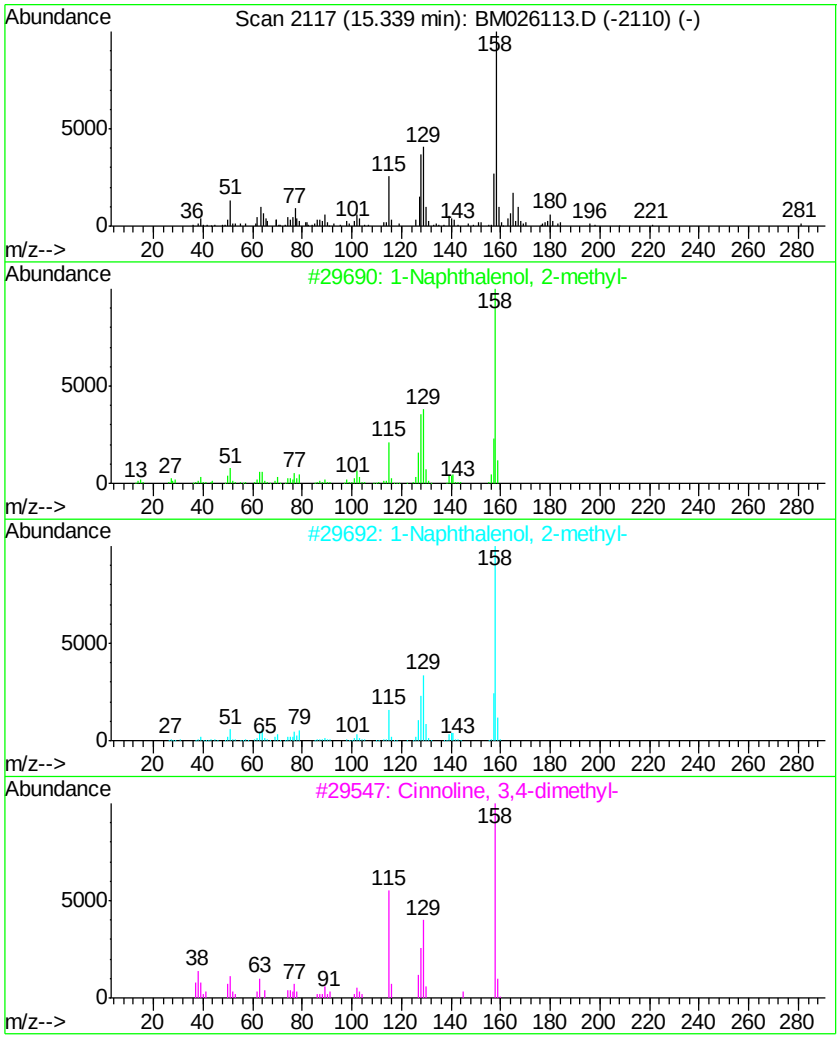
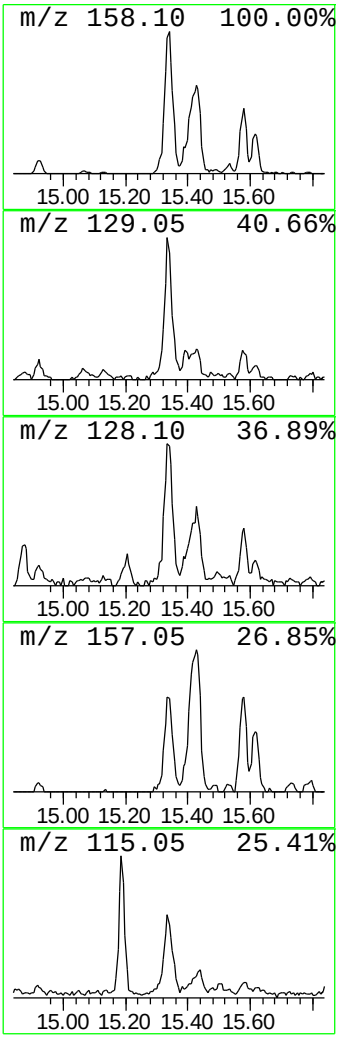
Peak Number 21 1-Naphthalenol, 2-methyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.34	3.49 ng/ul	280362	Acenaphthene-d10	14.13

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Naphthalenol, 2-methyl-	158	C11H10O	007469-77-4	93
2	1-Naphthalenol, 2-methyl-	158	C11H10O	007469-77-4	93
3	Cinnoline, 3,4-dimethyl-	158	C10H10N2	003929-83-7	70
4	7-Methyl-1-naphthol	158	C11H10O	006939-33-9	68
5	1H-Pyrazole, 3-methyl-5-phenyl-	158	C10H10N2	003347-62-4	49





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Data File : BM026113.D
Acq On : 26 May 2020 14:26
Operator : MA/SJ
Sample : L2737-16DL 10X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9B13DL

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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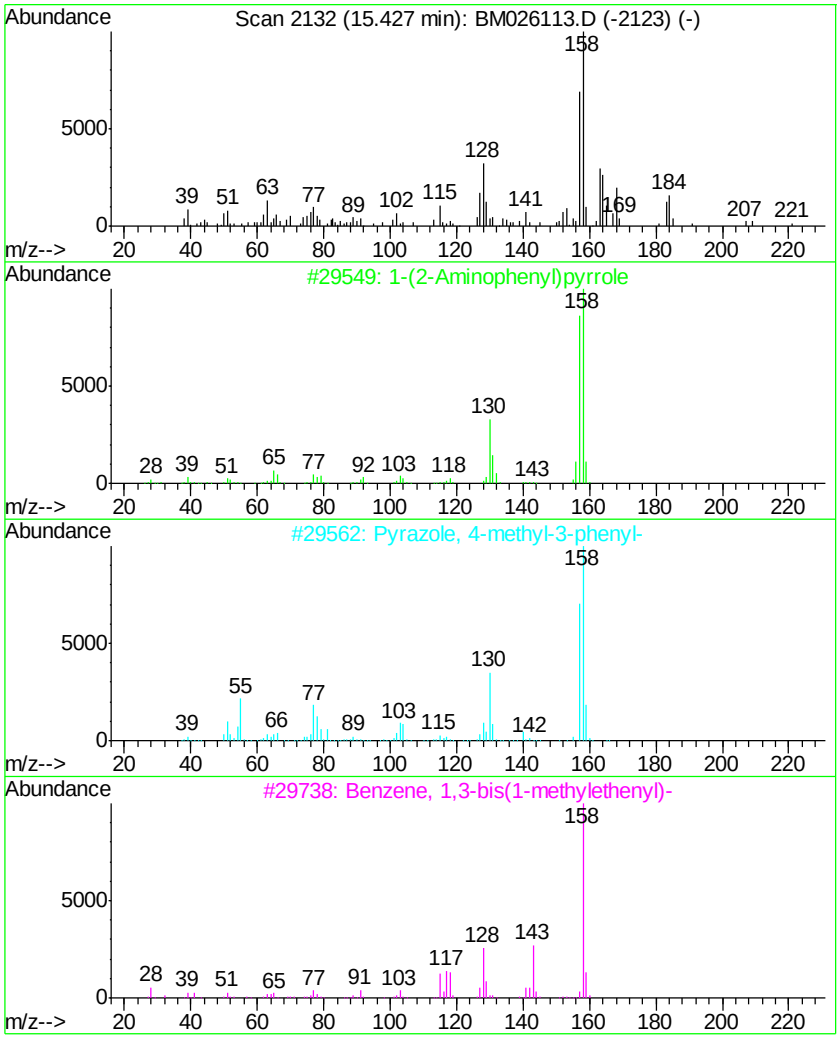
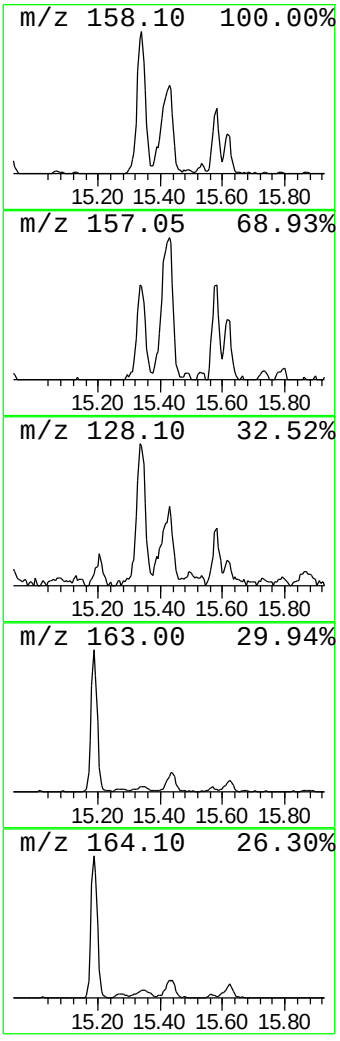
Peak Number 22 unknown-01 Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.43	2.79 ng/ul	224587	Acenaphthene-d10	14.13

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-(2-Aminophenyl)pyrrole	158	C10H10N2	006025-60-1	38
2	Pyrazole, 4-methyl-3-phenyl-	158	C10H10N2	013808-62-3	35
3	Benzene, 1,3-bis(1-methylethenyl)-	158	C12H14	003748-13-8	27
4	1,4-Naphthalenediamine	158	C10H10N2	002243-61-0	27
5	Cinnoline, 3,4-dimethyl-	158	C10H10N2	003929-83-7	27





Library Search Compound Report

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Data File : BM026113.D
Acq On : 26 May 2020 14:26
Operator : MA/SJ
Sample : L2737-16DL 10X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9B13DL

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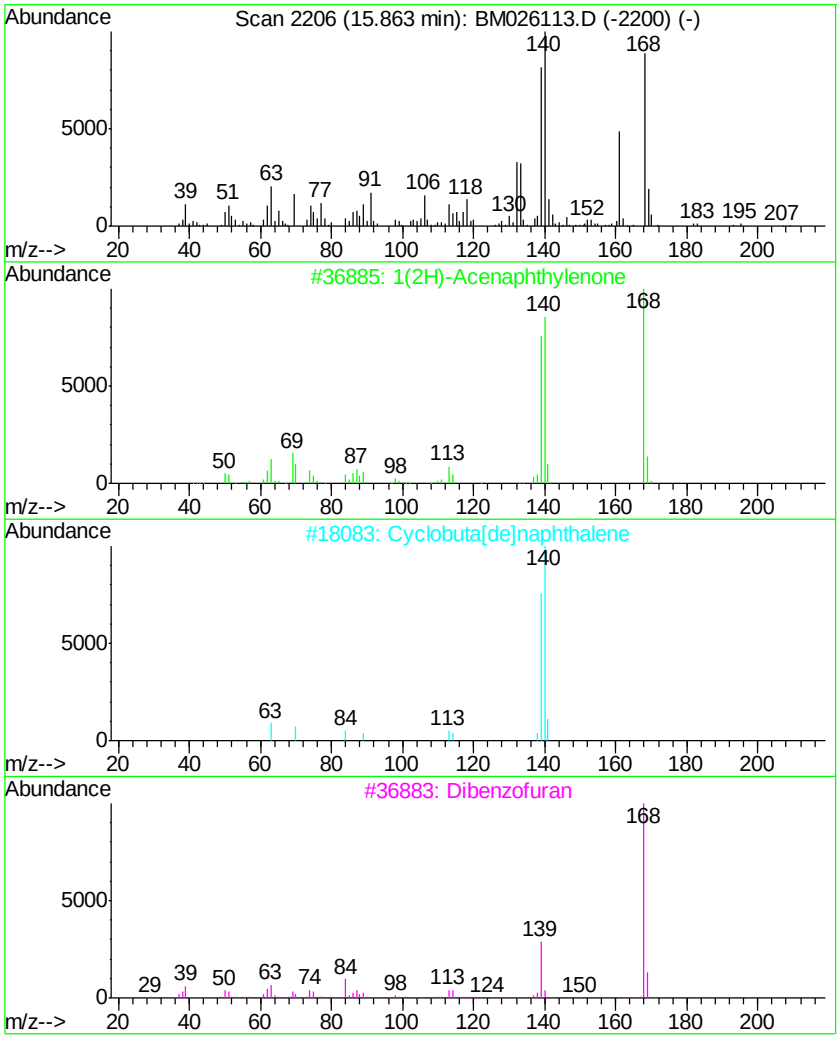
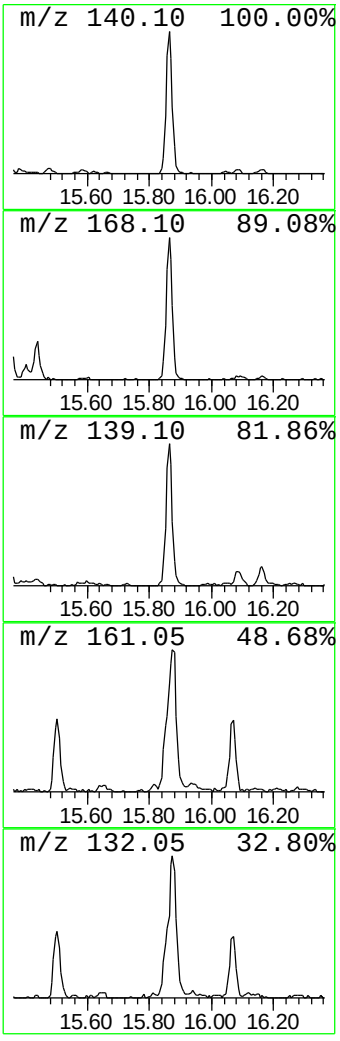
Peak Number 23 1(2H)-Acenaphthylenone Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.86	3.60 ng/ul	354101	Phenanthrene-d10	16.88

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1(2H)-Acenaphthylenone	168	C12H8O	002235-15-6	92
2	Cyclobuta[de]naphthalene	140	C11H8	024973-91-9	38
3	Dibenzofuran	168	C12H8O	000132-64-9	35
4	Benzaldehyde, 2-fluoro-4-hydroxy-	140	C7H5FO2	000348-27-6	32
5	2,5-Thiophenedicarboxaldehyde	140	C6H4O2S	000932-95-6	30





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Acq On : 26 May 2020 14:26
Operator : MA/SJ
Sample : L2737-16DL 10X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9B13DL

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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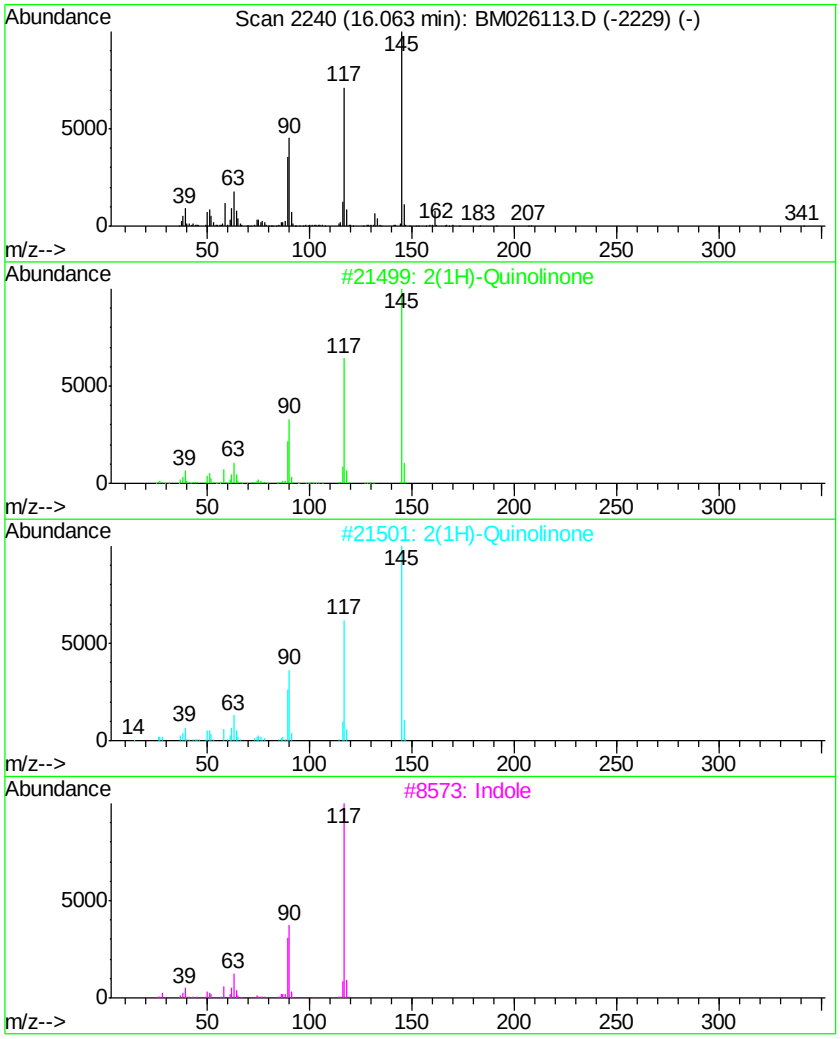
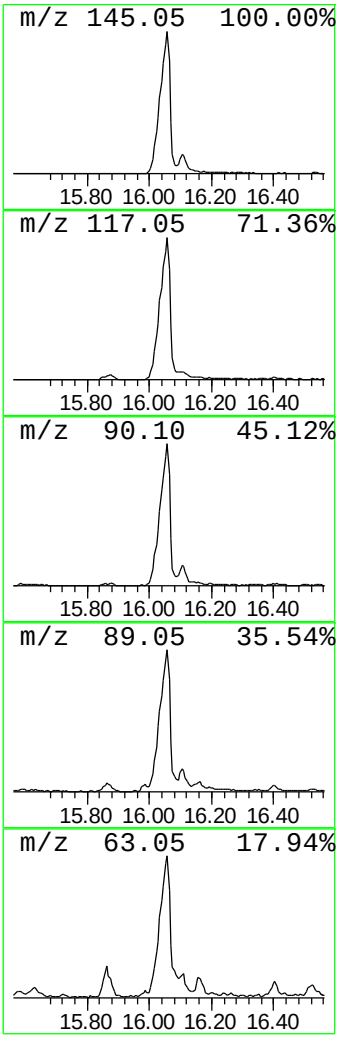
Peak Number 24 2(1H)-Quinolinone Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.06	11.04 ng/ul	1084370	Phenanthrene-d10	16.88

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	96
3	Indole	117	C8H7N	000120-72-9	95
4	8-Hydroxyquinoline	145	C9H7NO	000148-24-3	94
5	5-Quinolinol	145	C9H7NO	000578-67-6	94





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Acq On : 26 May 2020 14:26
Operator : MA/SJ
Sample : L2737-16DL 10X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9B13DL

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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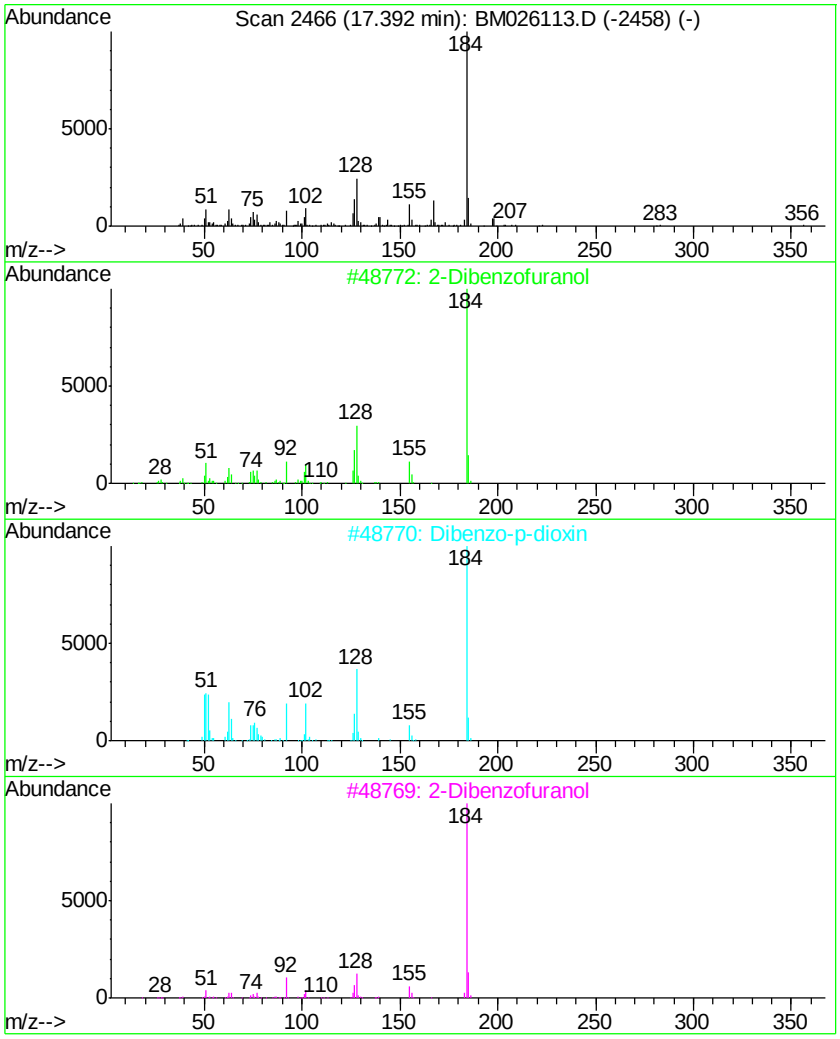
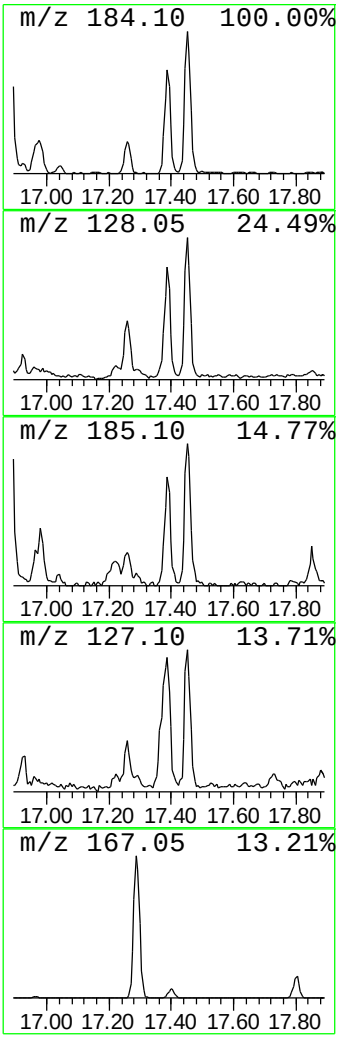
Peak Number 25 2-Dibenzofuranol Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.39	3.02 ng/ul	296738	Phenanthrene-d10	16.88

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Dibenzofuranol	184	C12H8O2	000086-77-1	98
2	Dibenzo-p-dioxin	184	C12H8O2	000262-12-4	90
3	2-Dibenzofuranol	184	C12H8O2	000086-77-1	90
4	Dibenzo-p-dioxin	184	C12H8O2	000262-12-4	87
5	Dibenzo-p-dioxin	184	C12H8O2	000262-12-4	87





Library Search Compound Report

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Sample : L2737-16DL 10X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9B13DL

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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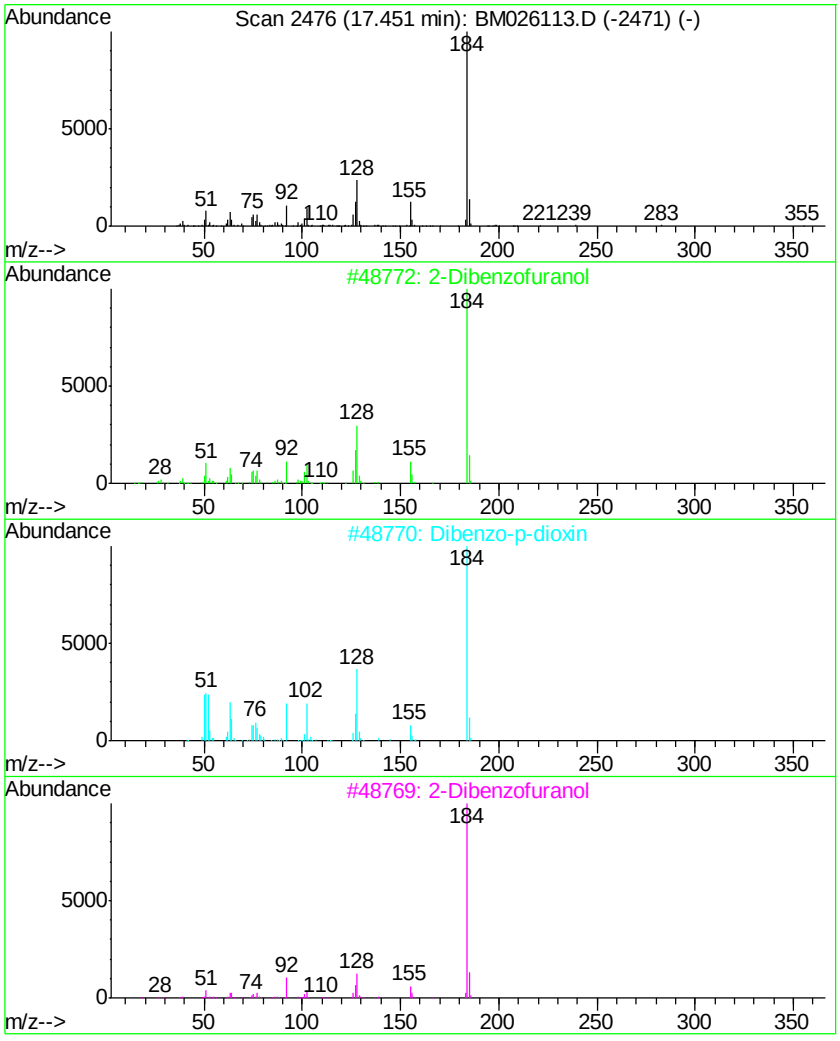
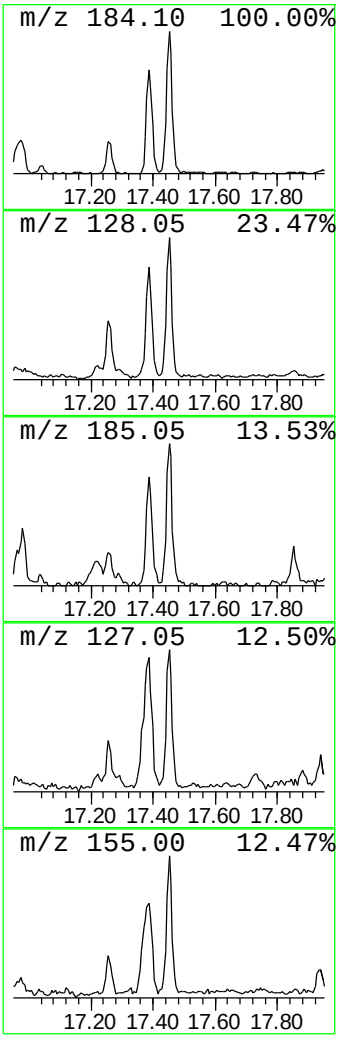
Peak Number 26 Dibenzo-p-dioxin Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.45	2.59 ng/ul	254040	Phenanthrene-d10	16.88

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Dibenzofuranol	184	C12H8O2	000086-77-1	97
2	Dibenzo-p-dioxin	184	C12H8O2	000262-12-4	94
3	2-Dibenzofuranol	184	C12H8O2	000086-77-1	93
4	2-Dibenzofuranol	184	C12H8O2	000086-77-1	86
5	Dibenzo-p-dioxin	184	C12H8O2	000262-12-4	86





Tentatively Identified Compound (LSC) summary

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Data File : BM026113.D
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Operator : MA/SJ
Sample : L2737-16DL 10X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F9B13DL

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	2.1	ng/ul	92872	1	7.49	902918	20.0
Indane	7.86	20.5	ng/ul	925098	1	7.49	902918	20.0
Benzene, 1-propynyl-	8.02	2.9	ng/ul	132443	1	7.49	902918	20.0
Benzonitrile, 2-m...	8.33	2.6	ng/ul	116055	1	7.49	902918	20.0
Phenol, 2,6-dimet...	8.90	4.5	ng/ul	272233	2	10.25	1222680	20.0
Cyclohexanemethan...	9.64	6.4	ng/ul	392085	2	10.25	1222680	20.0
Phenol, 4-ethyl-	9.71	14.7	ng/ul	897270	2	10.25	1222680	20.0
Phenol, 3,5-dimet...	9.76	43.2	ng/ul	2643140	2	10.25	1222680	20.0
Phenol, 3,4-dimet...	10.15	10.6	ng/ul	646503	2	10.25	1222680	20.0
Benzo[b]thiophene	10.42	15.2	ng/ul	926060	2	10.25	1222680	20.0
Phenol, 4-ethyl-3...	10.81	2.0	ng/ul	124125	2	10.25	1222680	20.0
Phenol, 2-propyl-	11.10	7.3	ng/ul	442923	2	10.25	1222680	20.0
Phenol, 2,4,6-tri...	11.35	2.2	ng/ul	136496	2	10.25	1222680	20.0
1H-Inden-5-ol, 2,...	12.03	5.2	ng/ul	315658	2	10.25	1222680	20.0
Naphthalene, 1-me...	12.15	17.6	ng/ul	1072670	2	10.25	1222680	20.0
Benzaldehyde, 3,4...	12.33	2.8	ng/ul	220943	3	14.13	1608930	20.0
Naphthalene, 2,3-...	13.45	2.4	ng/ul	190428	3	14.13	1608930	20.0
2-Naphthalenol	14.33	4.7	ng/ul	379108	3	14.13	1608930	20.0
Benzofuran, 2-eth...	14.43	7.1	ng/ul	572610	3	14.13	1608930	20.0
Pyridine, 5-ethen...	14.87	3.1	ng/ul	248510	3	14.13	1608930	20.0
1-Naphthalenol, 2...	15.34	3.5	ng/ul	280362	3	14.13	1608930	20.0
unknown-01	15.43	2.8	ng/ul	224587	3	14.13	1608930	20.0
1(2H)-Acenaphthyl...	15.86	3.6	ng/ul	354101	4	16.88	1964570	20.0
2(1H)-Quinolinone	16.06	11.0	ng/ul	1084370	4	16.88	1964570	20.0
2-Dibenzofuranol	17.39	3.0	ng/ul	296738	4	16.88	1964570	20.0
Dibenzo-p-dioxin	17.45	2.6	ng/ul	254040	4	16.88	1964570	20.0