LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100518\
Data File : BM017051.D
Acq On : 06 Oct 2018 00:41
Operator : MJ/SJ
Sample : J5298-06
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : E62T9

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-BM100418MA.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	3.963	163	166	172	rVB	95293	129035	1.80%	0.125%
2	6.869	654	660	673	rVB	310926	526915	7.35%	0.512%
3	7.040	683	689	699	rVB	1039620	1525879	21.29%	1.483%
4	7.228	715	721	732	rVB	1079326	1732549	24.17%	1.684%
5	7.698	794	801	807	rVB	978493	1557568	21.73%	1.514%
6	8.063	858	863	871	rBV2	108886	227790	3.18%	0.221%
7	8.228	885	891	898	rVB	74518	113121	1.58%	0.110%
8	8.398	914	920	928	rVB	889545	1394904	19.46%	1.356%
9	8.663	961	965	970	rVB2	128210	188417	2.63%	0.183%
10	8.763	977	982	986	rVV	184841	297477	4.15%	0.289%
11	8.851	990	997	1004	rVV	1265374	2136469	29.81%	2.077%
12	8.957	1010	1015	1020	rVB2	70659	107256	1.50%	0.104%
13	9.045	1020	1030	1037	rVB4	112073	306228	4.27%	0.298%
14	9.151	1043	1048	1055	rVV	178729	375671	5.24%	0.365%
15	9.222	1055	1060	1071	rVB	507323	886043	12.36%	0.861%
16	9.563	1111	1118	1124	rBV	1049342	1724065	24.06%	1.676%
17	9.610	1124	1126	1135	rVB	96485	127255	1.78%	0.124%
18	9.728	1142	1146	1151	rBV2	64285	95433	1.33%	0.093%
19	9.992	1182	1191	1200	rVB4	56653	158417	2.21%	0.154%
20	10.098	1200	1209	1219	rBV	1607367	2931433	40.90%	2.850%
21	10.304	1239	1244	1252	rBV	176859	342104	4.77%	0.333%
22	10.380	1253	1257	1261	rVB2	83991	131482	1.83%	0.128%
23	10.475	1261	1273	1279	rBV	1554316	2732434	38.13%	2.656%
24	10.539	1279	1284	1287	rBV5	68466	124170	1.73%	0.121%
25	10.622	1292	1298	1307	rVB	959792	1638331	22.86%	1.593%
26	10.722	1309	1315	1324	rVB4	98958	245841	3.43%	0.239%
27	10.886	1339	1343	1350	rVB	144679	242237	3.38%	0.235%
28	11.022	1358	1366	1368	rBV	182103	298155	4.16%	0.290%
29	11.063	1368	1373	1380	rVB	725708	1221735	17.05%	1.188%
30	11.239	1394	1403	1410	rVB3	94029	185979	2.59%	0.181%
31	11.351	1416	1422	1427	rBV2	278136	491057	6.85%	0.477%
32	11.404	1427	1431	1436	rVB2	119428	199278	2.78%	0.194%
33	11.539	1447	1454	1458	rBV	53586	112523	1.57%	0.109%
34	11.592	1458	1463	1473	rBV2	260341	499900	6.98%	0.486%

LSC Area Percent Report

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Operator : MJ/SJ
Sample : J5298-06
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : E62T9

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-BM100418MA.M
Title : SVOA CALIBRATION

35	11.680	1473	1478	1482	rBV	180441	255680	3.57%	0.249%
36	11.804	1495	1499	1505	rVB4	56778	108403	1.51%	0.105%
37	11.886	1509	1513	1518	rVB	201193	299989	4.19%	0.292%
38	12.027	1532	1537	1540	rBV3	71065	130841	1.83%	0.127%
39	12.186	1559	1564	1569	rBV	633440	920166	12.84%	0.895%
40	12.245	1569	1574	1582	rVB4	316967	656839	9.16%	0.639%
41	12.333	1582	1589	1592	rBV2	140255	292485	4.08%	0.284%
42	12.480	1608	1614	1618	rBV	375086	545652	7.61%	0.530%
43	12.527	1618	1622	1626	rVV2	193920	357042	4.98%	0.347%
44	12.592	1626	1633	1640	rVB2	333889	800242	11.17%	0.778%
45	12.686	1640	1649	1653	rBV4	665242	1342325	18.73%	1.305%
46	12.745	1653	1659	1665	rVB5	671769	1478255	20.63%	1.437%
47	12.810	1665	1670	1672	rBV	181418	289923	4.05%	0.282%
48	12.833	1672	1674	1685	rVB3	227903	402080	5.61%	0.391%
49	12.927	1685	1690	1697	rVB8	69361	140672	1.96%	0.137%
50	13.016	1697	1705	1706	rBV	118439	184910	2.58%	0.180%
51	13.074	1711	1715	1720	rVV3	158092	248736	3.47%	0.242%
52	13.127	1720	1724	1729	rVV2	173446	293273	4.09%	0.285%
53	13.204	1729	1737	1742	rVV9	206541	536133	7.48%	0.521%
54	13.304	1746	1754	1759	rVB4	384582	844112	11.78%	0.821%
55	13.469	1775	1782	1787	rBV2	365676	727544	10.15%	0.707%
56	13.527	1787	1792	1796	rVV4	286587	529794	7.39%	0.515%
57	13.574	1796	1800	1801	rVV	183341	237540	3.31%	0.231%
58	13.604	1801	1805	1809	rVV	377993	631875	8.82%	0.614%
59	13.686	1811	1819	1825	rVV4	473561	1635607	22.82%	1.590%
60	13.751	1825	1830	1836	rVB	3004411	4389916	61.25%	4.268%
61	13.839	1836	1845	1854	rBV3	902266	2090206	29.16%	2.032%
62	13.916	1854	1858	1863	rVV2	304507	655902	9.15%	0.638%
63	13.957	1863	1865	1868	rVV	195127	271269	3.79%	0.264%
64	14.027	1871	1877	1887	rVB	3691251	5495551	76.68%	5.342%
65	14.163	1896	1900	1903	rBV2	71213	90728	1.27%	0.088%
66	14.216	1903	1909	1914	rBV2	187217	413971	5.78%	0.402%
67	14.339	1923	1930	1939	rBV2	2261503	4115983	57.43%	4.001%
68	14.421	1939	1944	1948	rVV	396839	585171	8.16%	0.569%
69	14.510	1955	1959	1961	rBV2	255031	372317	5.19%	0.362%
70	14.539	1961	1964	1971	rVB2	728117	1031695	14.40%	1.003%
71	14.645	1977	1982	1985	rBV	154869	275934	3.85%	0.268%

LSC Area Percent Report

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Data File : BM017051.D
Acq On : 06 Oct 2018 00:41
Operator : MJ/SJ
Sample : J5298-06
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : E62T9

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-BM100418MA.M
Title : SVOA CALIBRATION

72	14.727	1992	1996	2000	rVB3	199822	303296	4.23%	0.295%
73	14.815	2007	2011	2018	rVB3	113162	225062	3.14%	0.219%
74	14.886	2018	2023	2027	rBV2	107872	207730	2.90%	0.202%
75	15.127	2060	2064	2069	rBV2	124201	188129	2.62%	0.183%
76	15.227	2078	2081	2085	rBV3	77819	96848	1.35%	0.094%
77	15.268	2085	2088	2091	rVV5	76978	127722	1.78%	0.124%
78	15.333	2093	2099	2105	rVV	4768351	6671198	93.08%	6.485%
79	15.457	2114	2120	2126	rVB2	1205410	1797630	25.08%	1.748%
80	15.557	2133	2137	2139	rBV5	68228	102619	1.43%	0.100%
81	15.586	2139	2142	2150	rVB2	145208	246674	3.44%	0.240%
82	15.668	2151	2156	2162	rVB8	40065	94582	1.32%	0.092%
83	15.768	2168	2173	2185	rBV4	123703	319637	4.46%	0.311%
84	15.962	2201	2206	2212	rBV7	69737	162755	2.27%	0.158%
85	16.086	2225	2227	2237	rVB6	89720	185074	2.58%	0.180%
86	16.210	2245	2248	2254	rVB2	75312	117144	1.63%	0.114%
87	16.392	2275	2279	2282	rBV3	134828	205439	2.87%	0.200%
88	16.539	2300	2304	2306	rBV4	76549	120628	1.68%	0.117%
89	16.651	2318	2323	2332	rBV7	81697	213631	2.98%	0.208%
90	16.762	2338	2342	2348	rVV4	65951	112967	1.58%	0.110%
91	16.821	2349	2352	2357	rVV4	85069	108875	1.52%	0.106%
92	17.004	2379	2383	2388	rVB2	116671	167668	2.34%	0.163%
93	17.086	2391	2397	2402	rVB	2743539	3832327	53.47%	3.726%
94	17.180	2408	2413	2426	rVB2	4596576	6724813	93.83%	6.537%
95	17.286	2427	2431	2439	rVB2	233265	388096	5.42%	0.377%
96	19.115	2737	2742	2747	rVB	81302	120845	1.69%	0.117%
97	19.480	2797	2804	2810	rBV	5433841	7166936	100.00%	6.967%
98	21.268	3103	3108	3115	rBV	2964197	3874107	54.06%	3.766%
99	23.356	3455	3463	3475	rBV	3433680	6740742	94.05%	6.553%
100	23.491	3480	3486	3496	rVB	1879650	3557944	49.64%	3.459%

Sum of corrected areas: 102867030

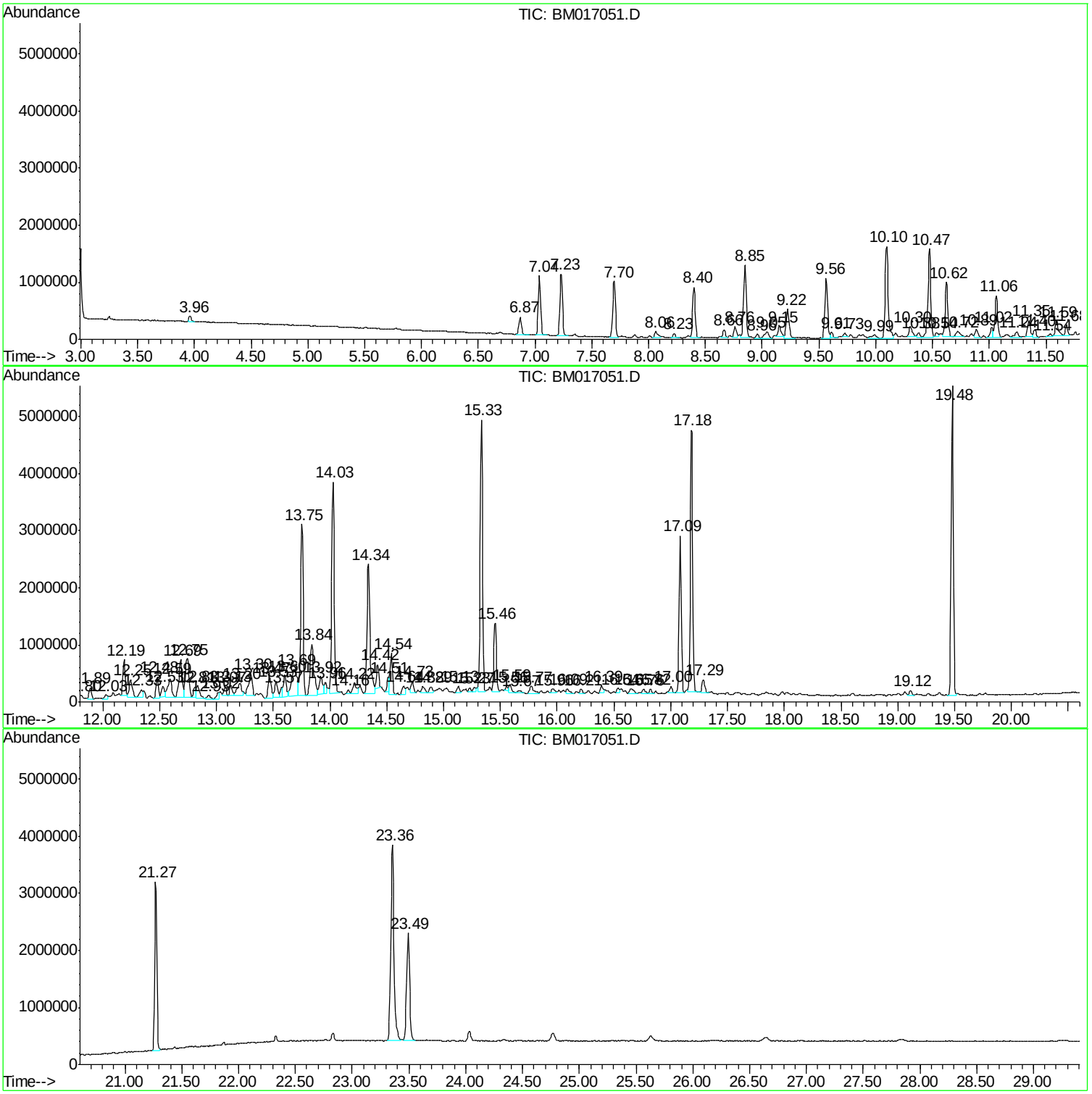
LSC Report - Integrated Chromatogram

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100518\
Data File : BM017051.D
Acq On : 06 Oct 2018 00:41
Operator : MJ/SJ
Sample : J5298-06
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : E62T9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM100418MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100518\
Data File : BM017051.D
Acq On : 06 Oct 2018 00:41
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Instrument : BNA_M
ClientSampleId : E62T9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM100418MA.M
Quant Title : SVOA CALIBRATION

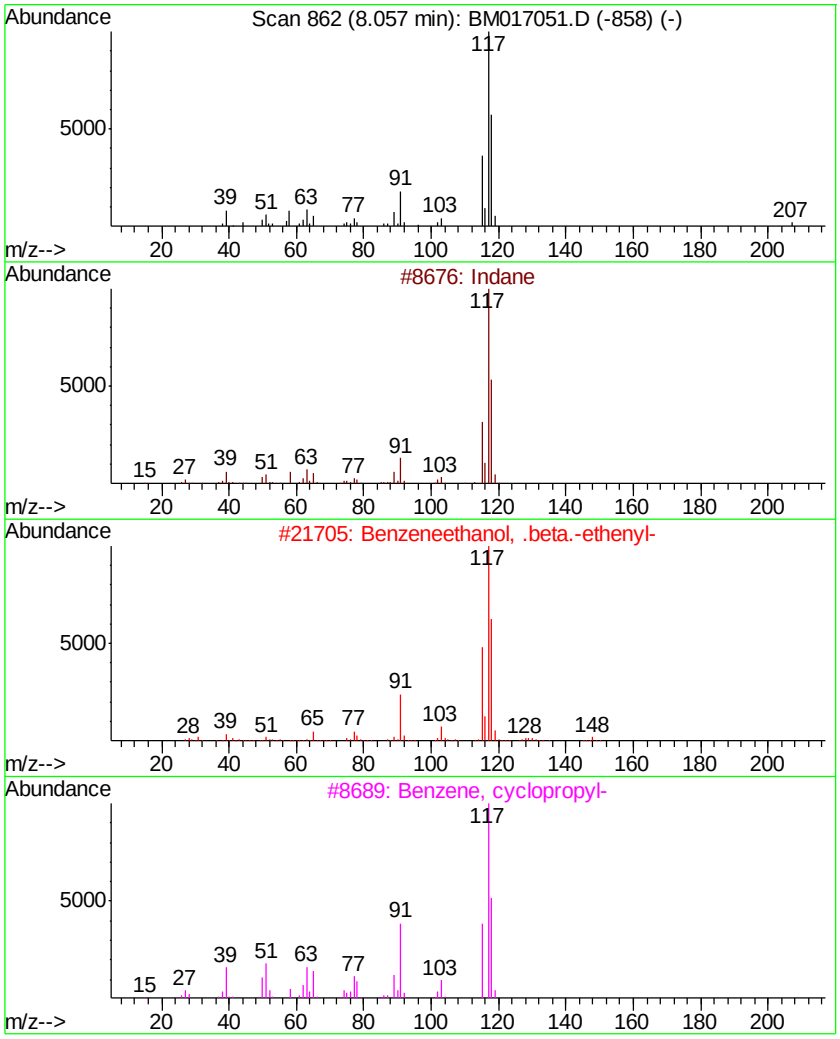
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

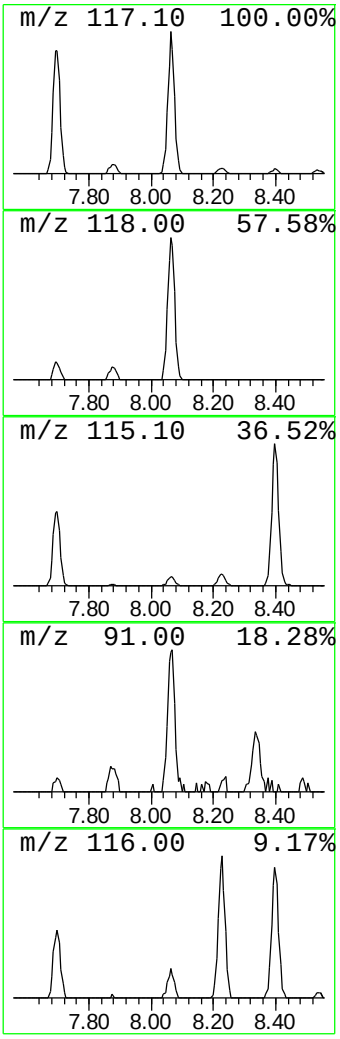
****

Peak Number 1 Indane Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.06	2.92 ng/ul	227790	1,4-Dichlorobenzene-d4	7.70

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	95
2	Benzeneethanol, .beta.-ethenyl-	148	C10H12O	006052-63-7	83
3	Benzene, cyclopropyl-	118	C9H10	000873-49-4	76
4	Tetracyclo[3.3.1.0(2,8).0(4,6)]-...	118	C9H10	1000191-13-7	72
5	Deltacyclene	118	C9H10	007785-10-6	72





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100518\
Data File : BM017051.D
Acq On : 06 Oct 2018 00:41
Operator : MJ/SJ
Sample : J5298-06
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : E62T9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM100418MA.M
Quant Title : SVOA CALIBRATION

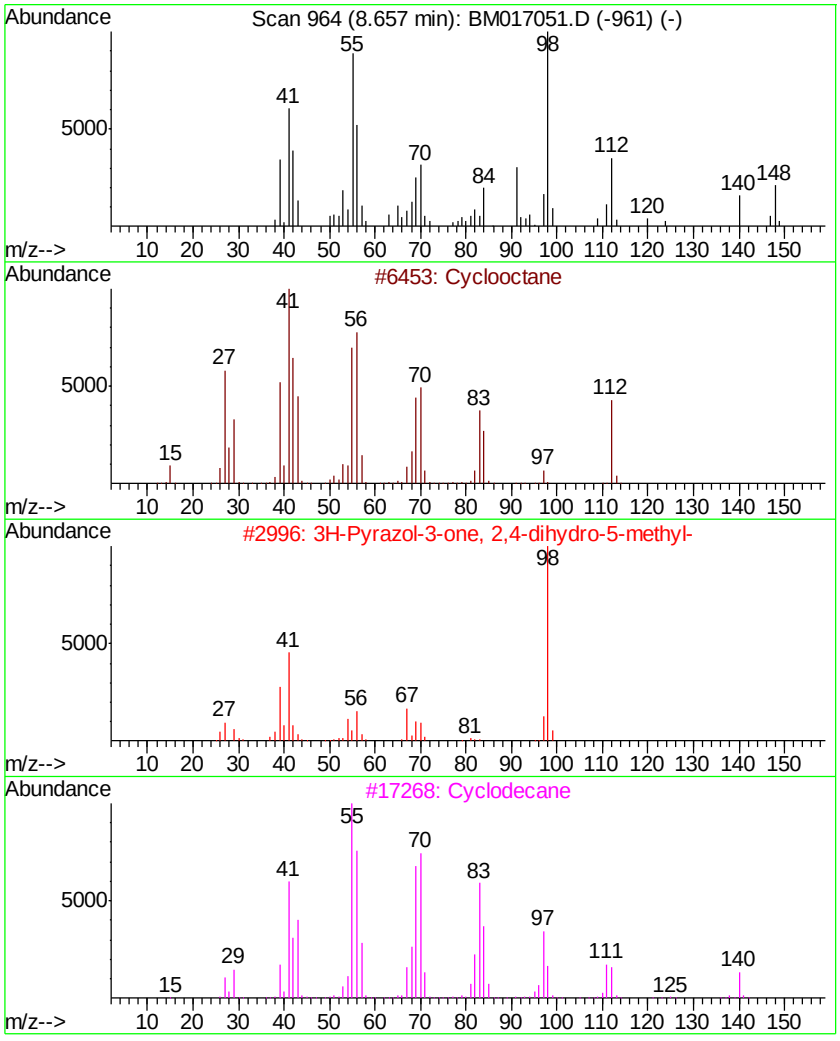
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

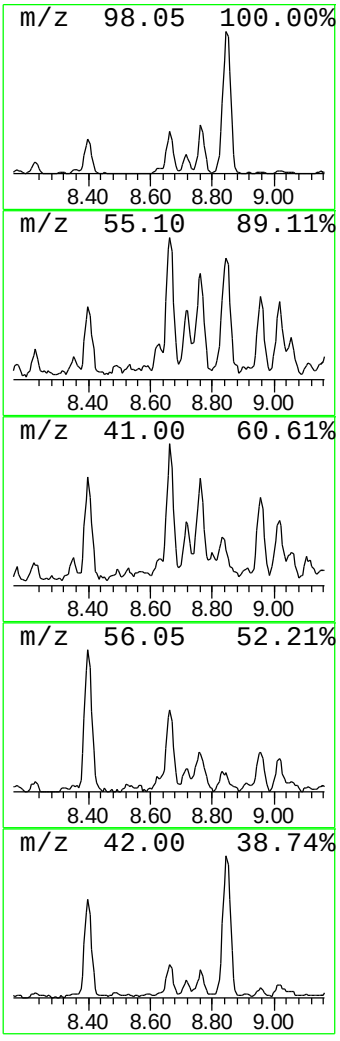
****

Peak Number 2 (DEL) Alkane: Cyclic8.66 Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.66	2.42 ng/ul	188417	1,4-Dichlorobenzene-d4	7.70

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclooctane	112	C8H16	000292-64-8	56
2	3H-Pyrazol-3-one, 2,4-dihydro-5-...	98	C4H6N2O	000108-26-9	52
3	Cyclodecane	140	C10H20	000293-96-9	46
4	Cyclooctane	112	C8H16	000292-64-8	38
5	cis-4-Decene	140	C10H20	019398-88-0	38





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100518\
Data File : BM017051.D
Acq On : 06 Oct 2018 00:41
Operator : MJ/SJ
Sample : J5298-06
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : E62T9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM100418MA.M
Quant Title : SVOA CALIBRATION

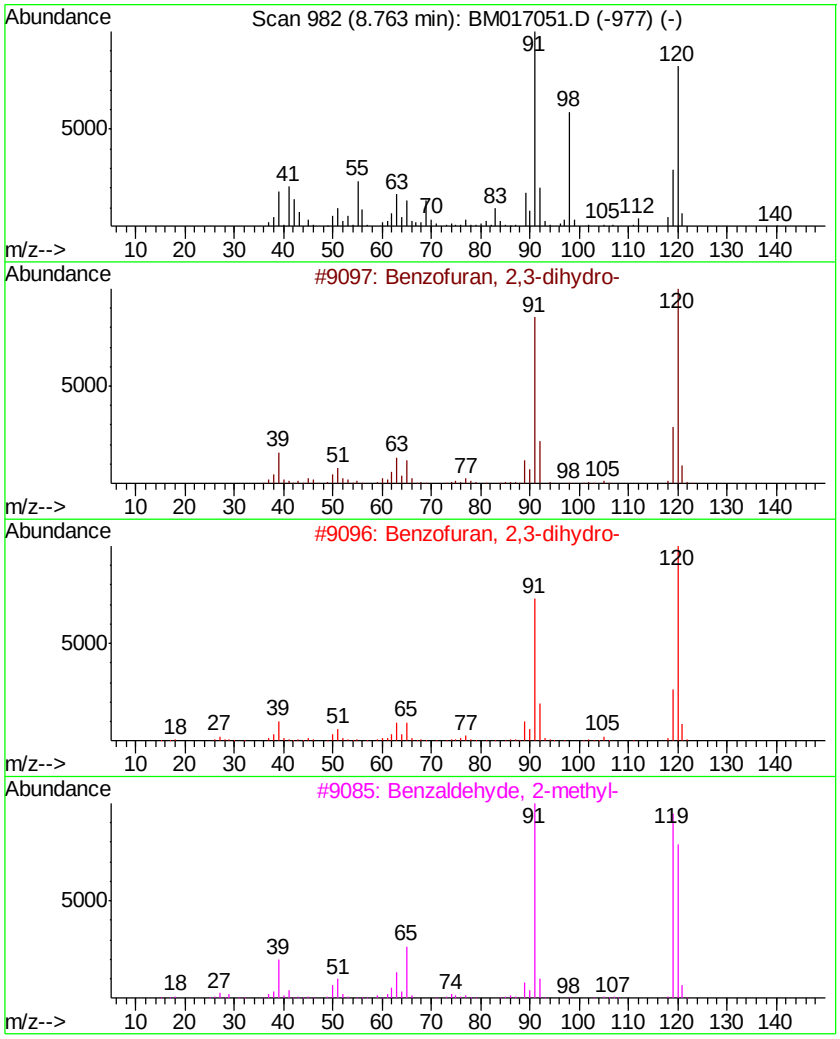
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

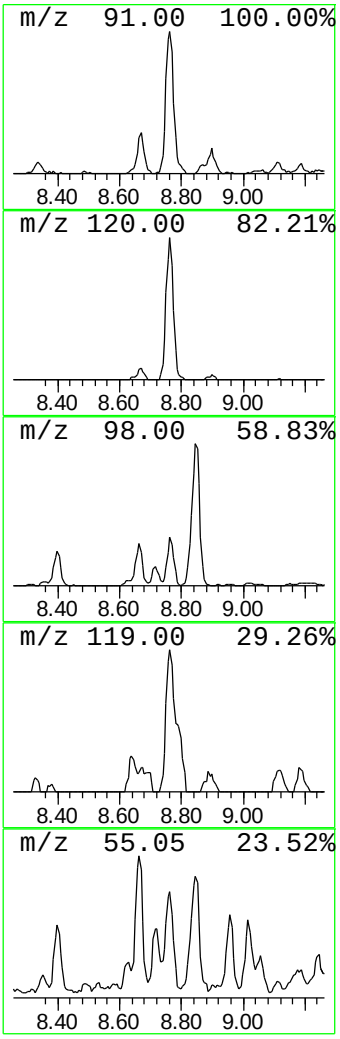
****

Peak Number 3 Benzofuran, 2,3-dihydro- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.76	3.82 ng/ul	297477	1,4-Dichlorobenzene-d4	7.70

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzofuran, 2,3-dihydro-	120	C8H8O	000496-16-2	86
2	Benzofuran, 2,3-dihydro-	120	C8H8O	000496-16-2	70
3	Benzaldehyde, 2-methyl-	120	C8H8O	000529-20-4	55
4	6-Methylenebicyclo[3.2.0]hept-3-...	120	C8H8O	1000211-11-9	49
5	Catecholborane	120	C6H5BO2	000274-07-7	49





Library Search Compound Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100518\
Data File : BM017051.D
Acq On : 06 Oct 2018 00:41
Operator : MJ/SJ
Sample : J5298-06
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : E62T9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM100418MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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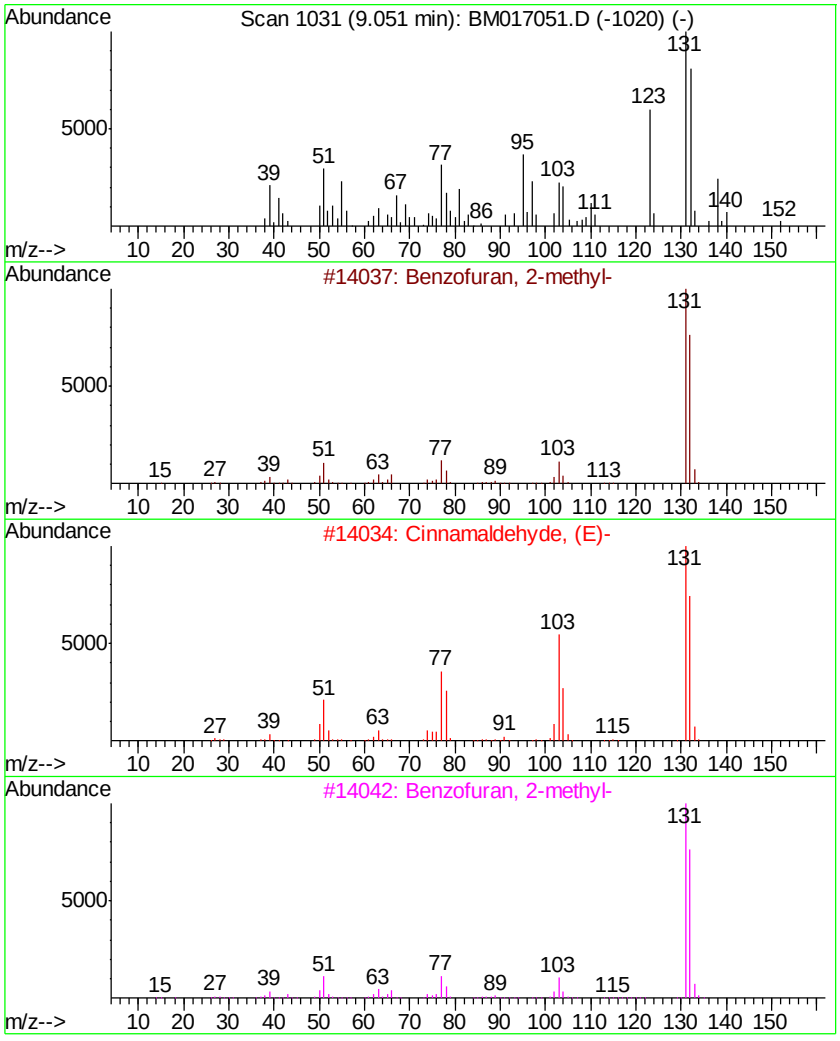
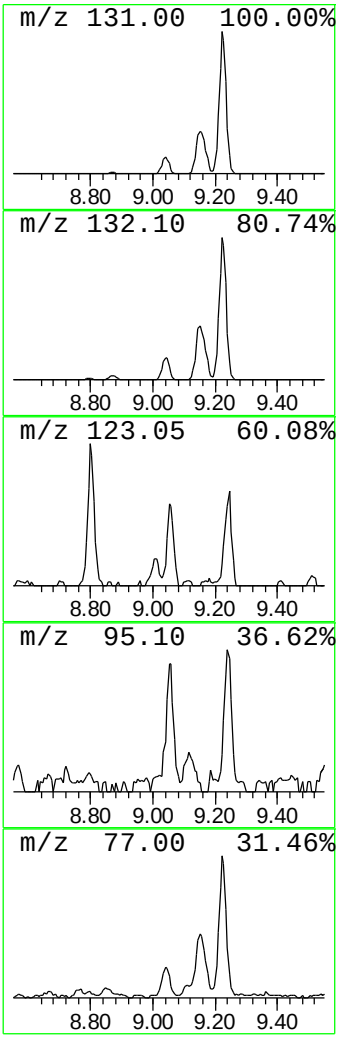
Peak Number 4 Benzofuran, 2-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.05	3.93 ng/ul	306228	1,4-Dichlorobenzene-d4	7.70

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	89
2	Cinnamaldehyde, (E)-	132	C9H8O	014371-10-9	81
3	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	76
4	Benzenemethanol, .alpha.-ethynyl-	132	C9H8O	004187-87-5	74
5	Benzofuran, 7-methyl-	132	C9H8O	017059-52-8	72





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100518\
Data File : BM017051.D
Acq On : 06 Oct 2018 00:41
Operator : MJ/SJ
Sample : J5298-06
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : E62T9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM100418MA.M
Quant Title : SVOA CALIBRATION

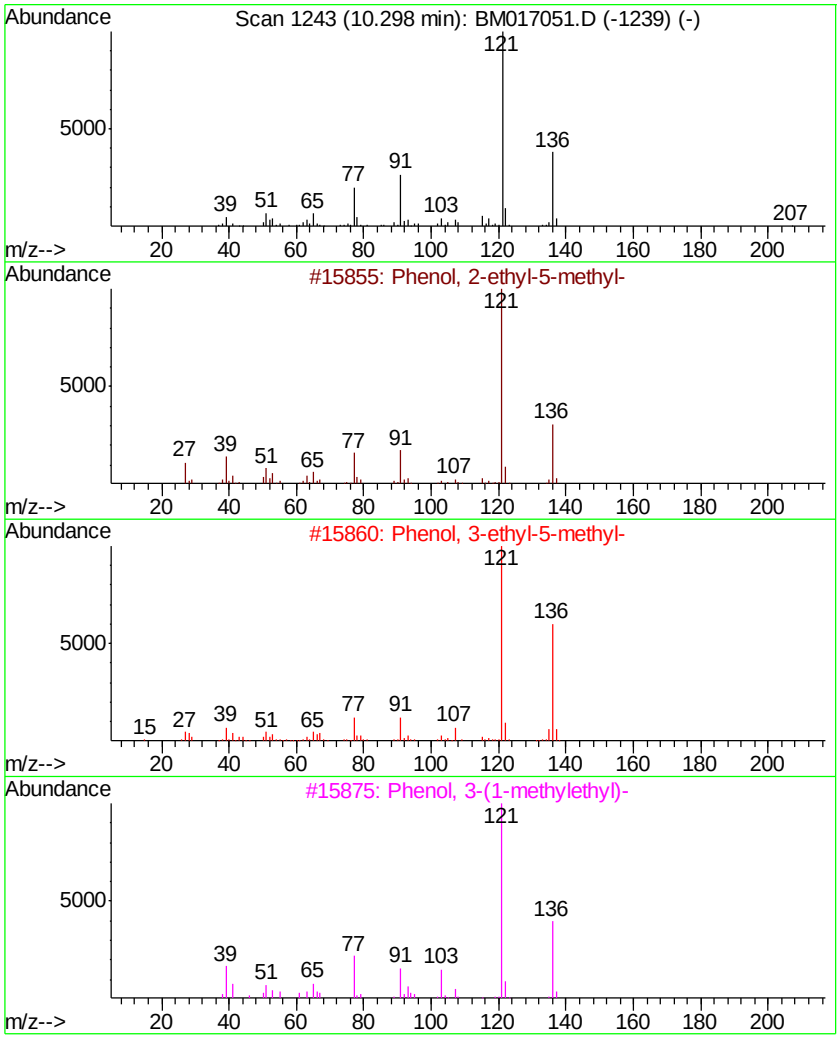
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

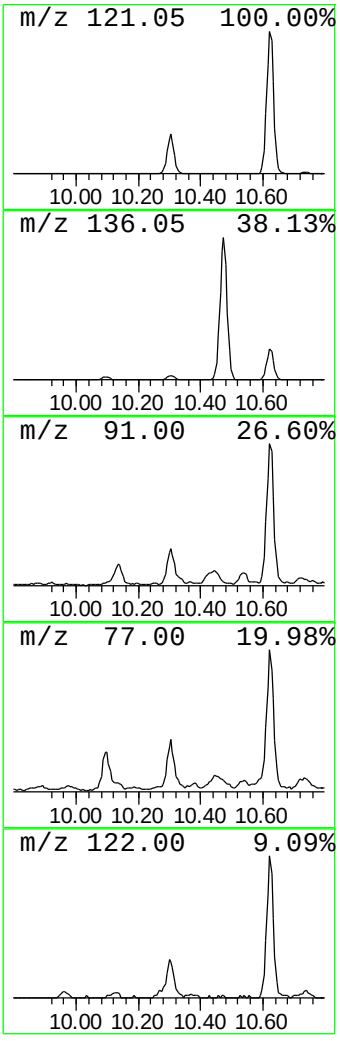
****

Peak Number 7 Phenol, 2-ethyl-5-methyl- Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.30	2.50 ng/ul	342104	Naphthalene-d8	10.47

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 2-ethyl-5-methyl-	136	C9H12O	001687-61-2	91
2	Phenol, 3-ethyl-5-methyl-	136	C9H12O	000698-71-5	90
3	Phenol, 3-(1-methylethyl)-	136	C9H12O	000618-45-1	86
4	Phenol, 4-ethyl-3-methyl-	136	C9H12O	001123-94-0	86
5	Phenol, 3-(1-methylethyl)-	136	C9H12O	000618-45-1	78





Library Search Compound Report

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Acq On : 06 Oct 2018 00:41
Operator : MJ/SJ
Sample : J5298-06
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : E62T9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM100418MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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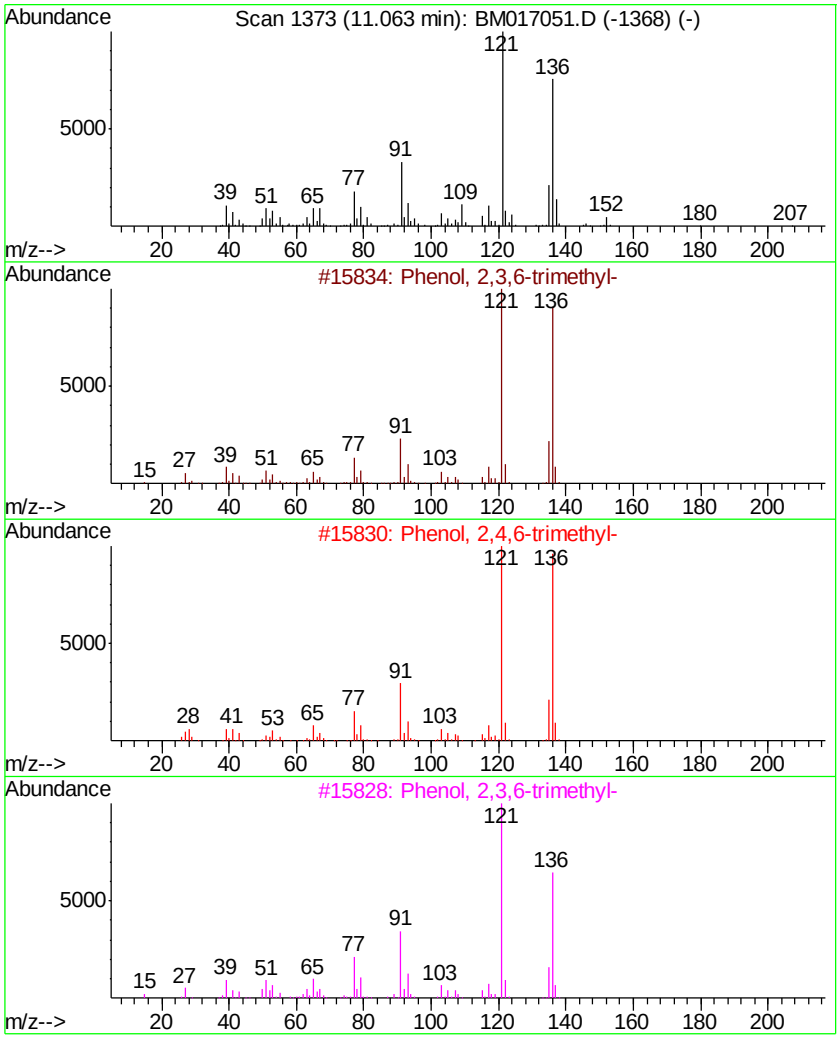
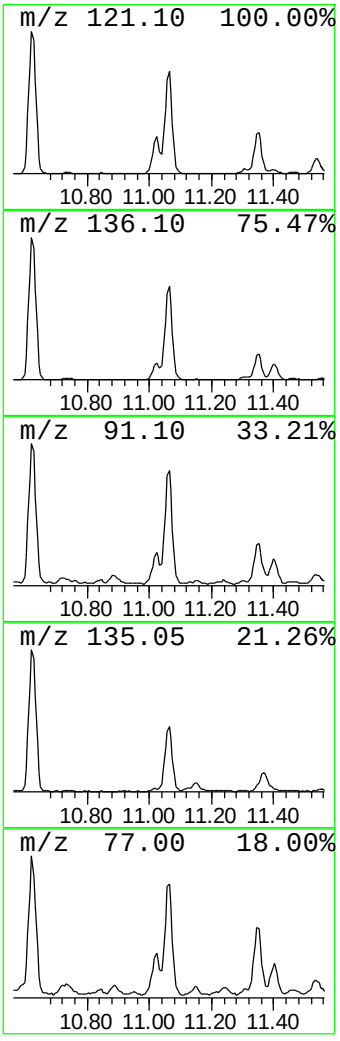
Peak Number 9 Phenol, 2,3,6-trimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.06	8.94 ng/ul	1221740	Naphthalene-d8	10.47

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





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Sample : J5298-06
Misc :
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Instrument : BNA_M
ClientSampleId : E62T9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM100418MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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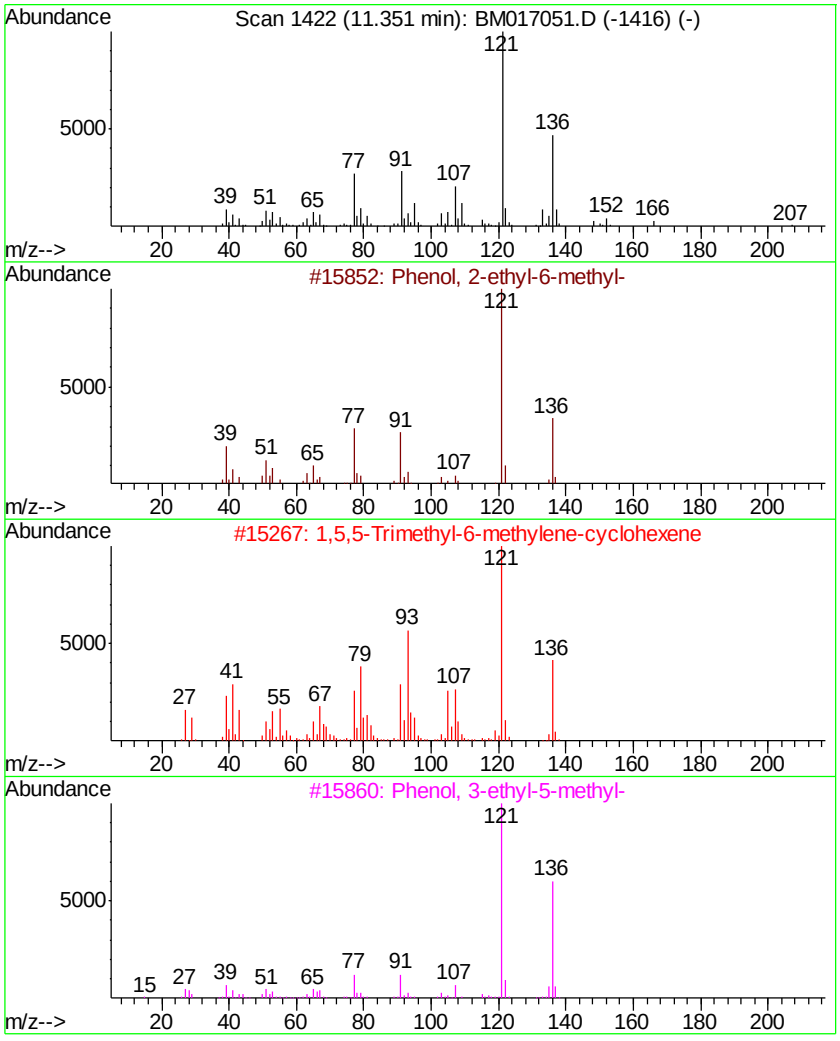
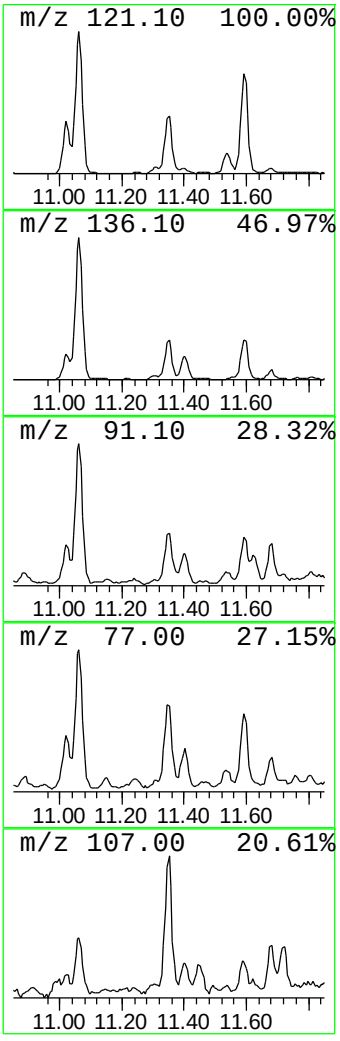
Peak Number 10 Phenol, 2-ethyl-6-methyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.35	3.59 ng/ul	491057	Naphthalene-d8	10.47

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 2-ethyl-6-methyl-	136	C9H12O	001687-64-5	81
2	1,5,5-Trimethyl-6-methylene-cycl...	136	C10H16	000514-95-4	80
3	Phenol, 3-ethyl-5-methyl-	136	C9H12O	000698-71-5	72
4	Phenol, 2-ethyl-4-methyl-	136	C9H12O	003855-26-3	72
5	Phenol, 2-(1-methylethyl)-	136	C9H12O	000088-69-7	68





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Sample : J5298-06
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : E62T9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM100418MA.M
Quant Title : SVOA CALIBRATION

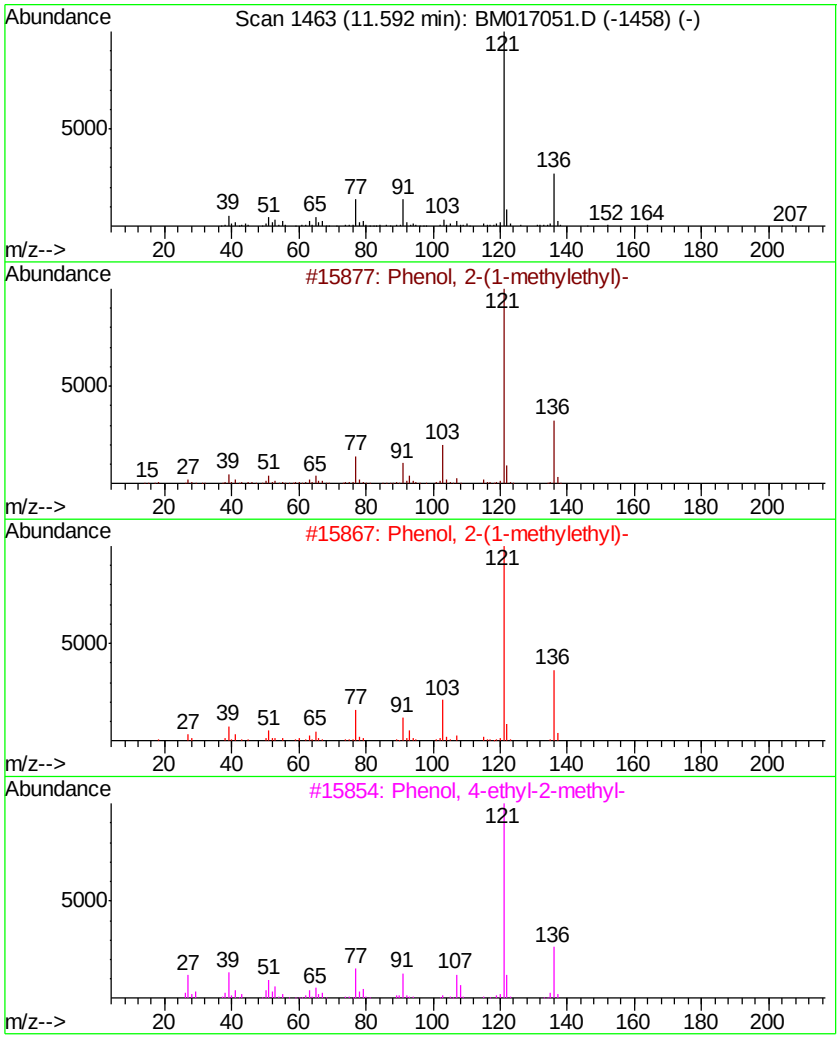
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

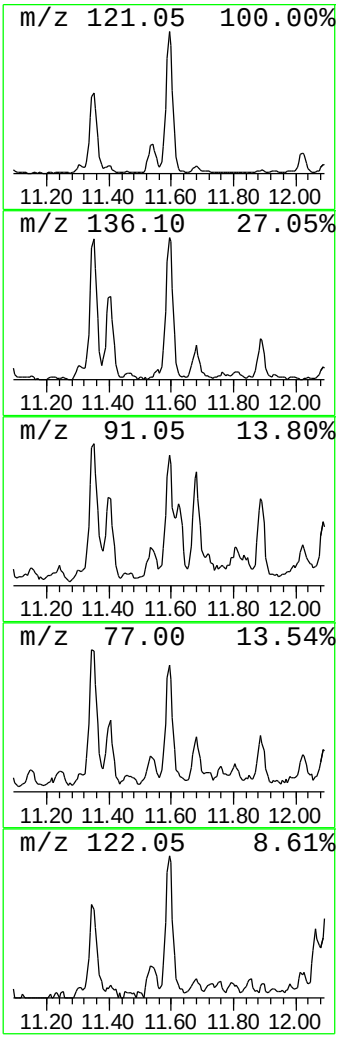
****

Peak Number 11 Phenol, 2-(1-methylethyl)- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.59	3.66 ng/ul	499900	Naphthalene-d8	10.47

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 2-(1-methylethyl)-	136	C9H12O	000088-69-7	91
2	Phenol, 2-(1-methylethyl)-	136	C9H12O	000088-69-7	91
3	Phenol, 4-ethyl-2-methyl-	136	C9H12O	002219-73-0	90
4	Phenol, 3-ethyl-5-methyl-	136	C9H12O	000698-71-5	90
5	Phenol, 4-(1-methylethyl)-	136	C9H12O	000099-89-8	90





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Operator : MJ/SJ
Sample : J5298-06
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : E62T9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM100418MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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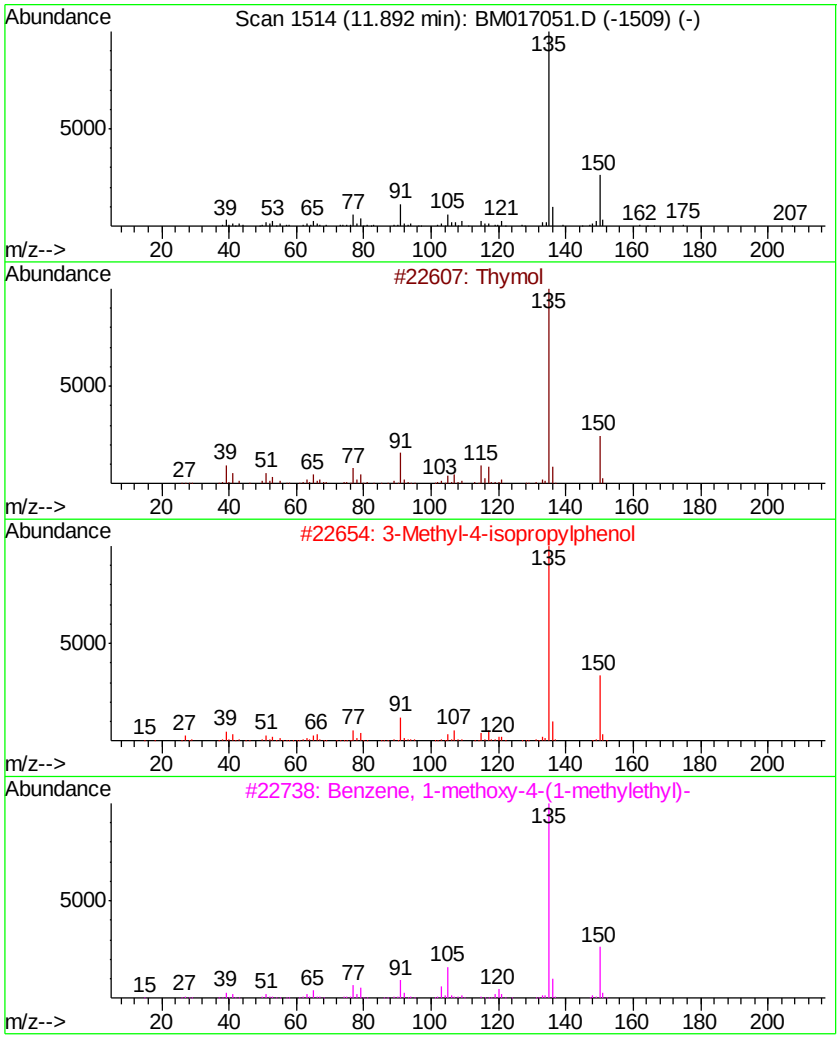
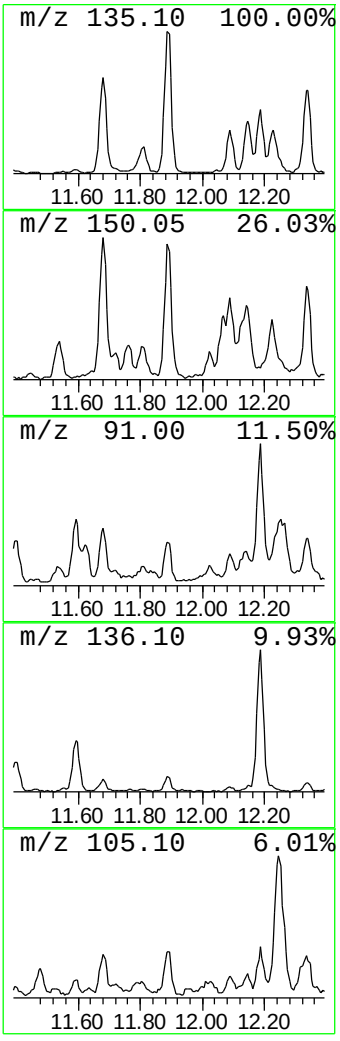
Peak Number 12 Thymol Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.89	2.20 ng/ul	299989	Naphthalene-d8	10.47

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Thymol	150	C10H14O	000089-83-8	91
2	3-Methyl-4-isopropylphenol	150	C10H14O	003228-02-2	91
3	Benzene, 1-methoxy-4-(1-methylet...	150	C10H14O	004132-48-3	91
4	Thymol	150	C10H14O	000089-83-8	91
5	Phenol, 2-methyl-5-(1-methylethyl)-	150	C10H14O	000499-75-2	90





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Acq On : 06 Oct 2018 00:41
Operator : MJ/SJ
Sample : J5298-06
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : E62T9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM100418MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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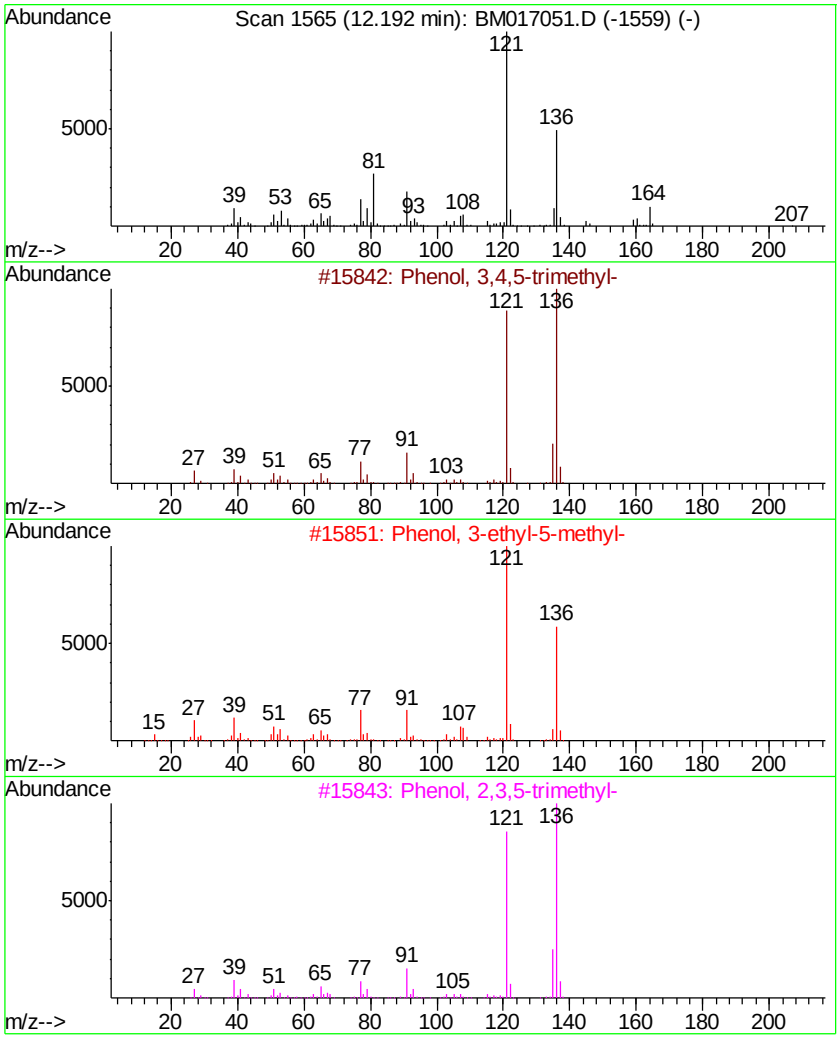
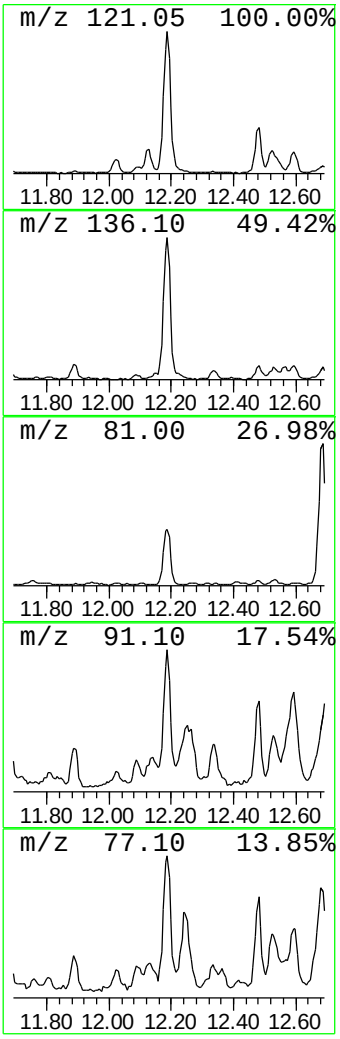
Peak Number 13 Phenol, 3,4,5-trimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.19	6.74 ng/ul	920166	Naphthalene-d8	10.47

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 3,4,5-trimethyl-	136	C9H12O	000527-54-8	90
2	Phenol, 3-ethyl-5-methyl-	136	C9H12O	000698-71-5	87
3	Phenol, 2,3,5-trimethyl-	136	C9H12O	000697-82-5	87
4	Phenol, 2-ethyl-4-methyl-	136	C9H12O	003855-26-3	83
5	Phenol, 3,4,5-trimethyl-	136	C9H12O	000527-54-8	83





Library Search Compound Report

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Operator : MJ/SJ
Sample : J5298-06
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : E62T9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM100418MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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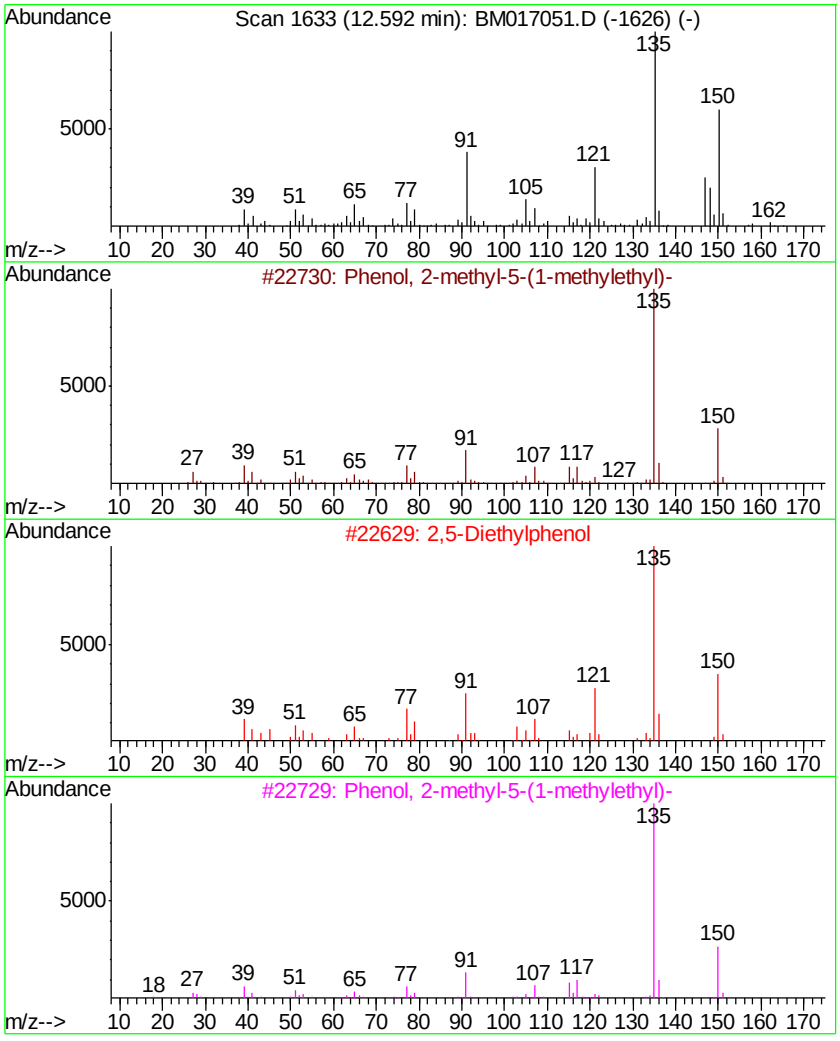
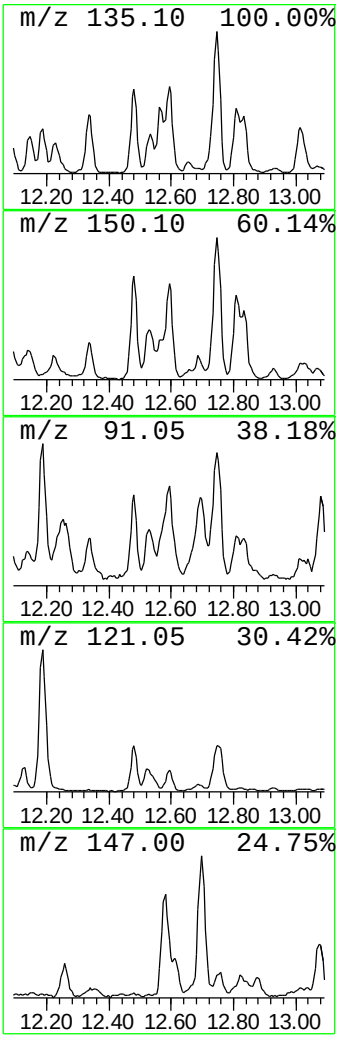
Peak Number 17 Phenol, 2-methyl-5-(1-methy... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.59	3.89 ng/ul	800242	Acenaphthene-d10	14.33

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 2-methyl-5-(1-methylethyl)-	150	C10H14O	000499-75-2	70
2	2,5-Diethylphenol	150	C10H14O	000876-20-0	70
3	Phenol, 2-methyl-5-(1-methylethyl)-	150	C10H14O	000499-75-2	70
4	Phenol, 2,3,4,6-tetramethyl-	150	C10H14O	003238-38-8	70
5	Phenol, 2-ethyl-4,5-dimethyl-	150	C10H14O	002219-78-5	62





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Sample : J5298-06
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : E62T9

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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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Peak Number 18 unknown-01 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.69	6.52 ng/ul	1342330	Acenaphthene-d10	14.33

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Isomyocorene	136	C10H16	006876-07-9	43
2	Cyclohexene, 3-(2-propynyl)-	120	C9H12	055956-43-9	41
3	1,4-Hexadiene, 3-ethyl-	110	C8H14	002080-89-9	38
4	Cyclohexene,3-(2-propenyl)-	122	C9H14	015232-95-8	38
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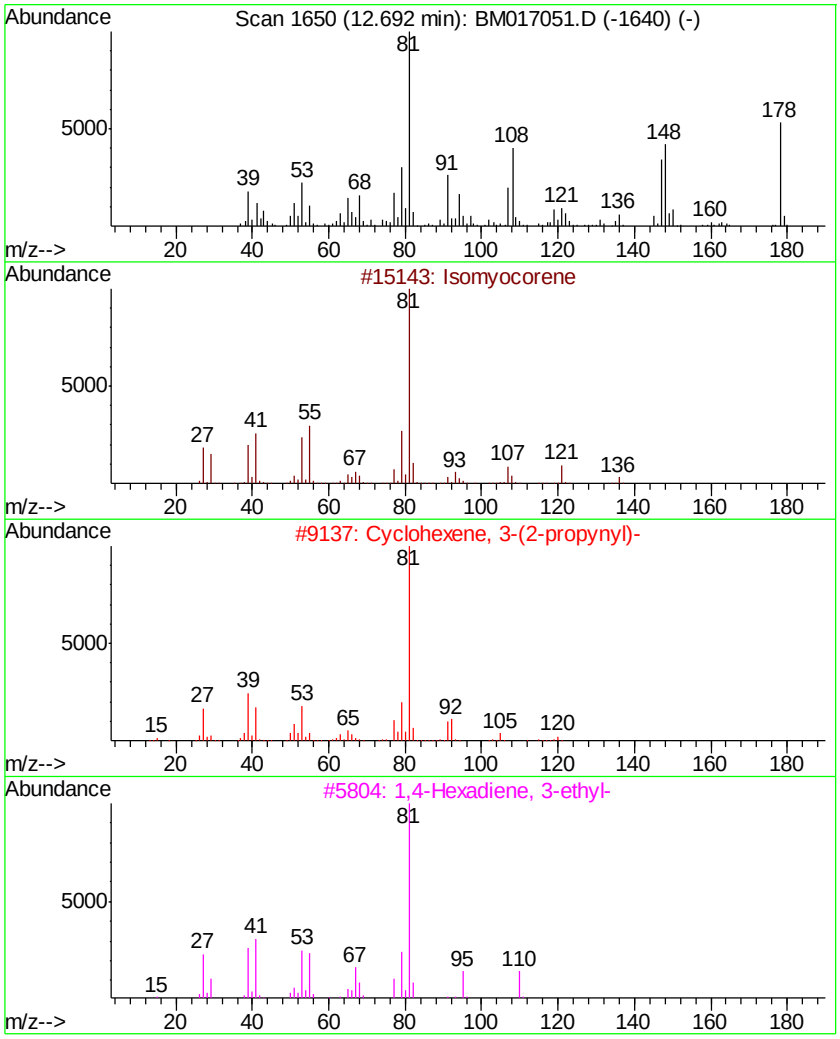

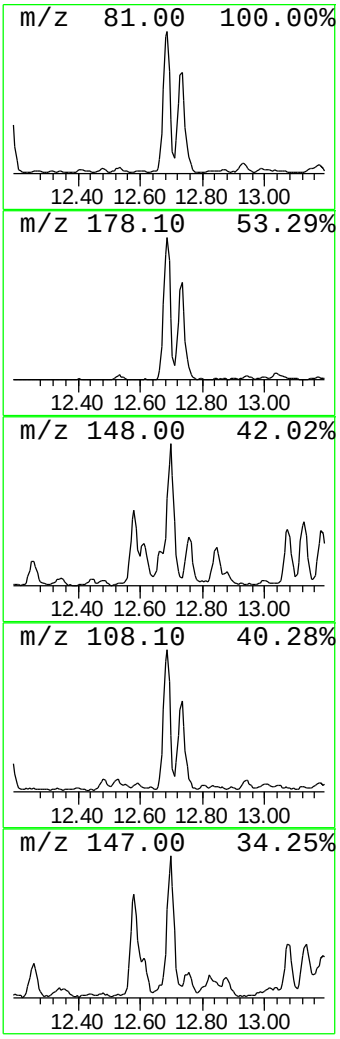

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Sample : J5298-06
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : E62T9

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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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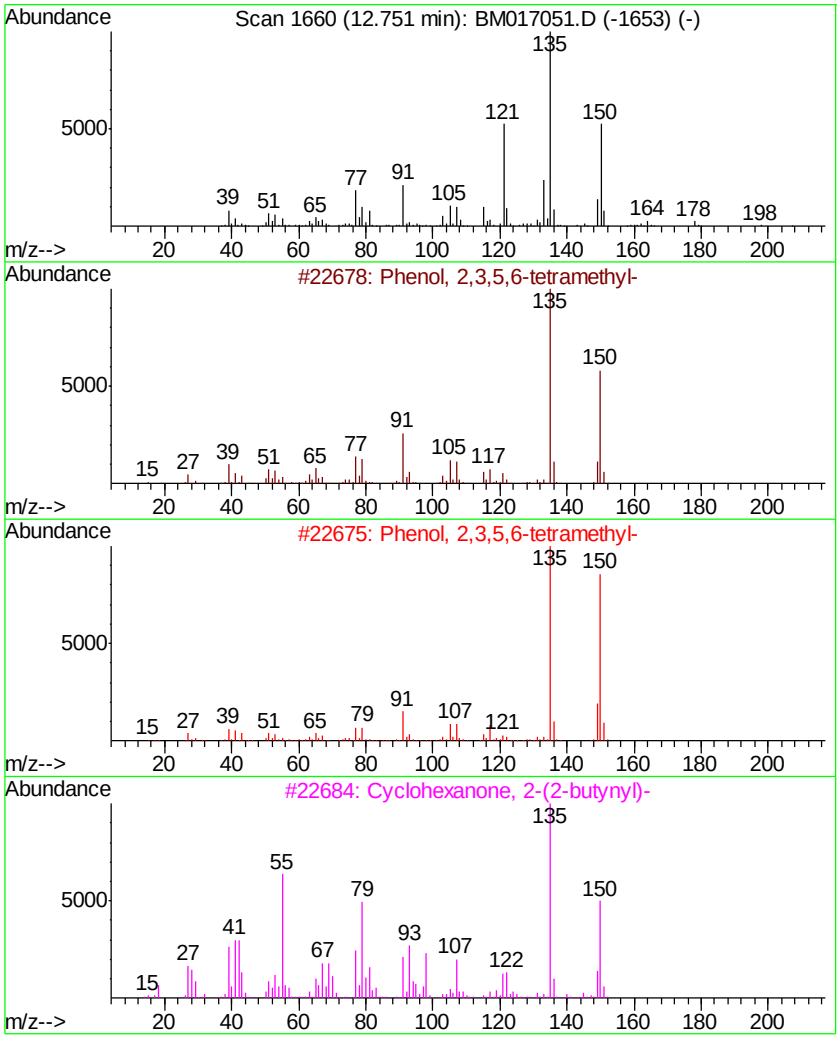
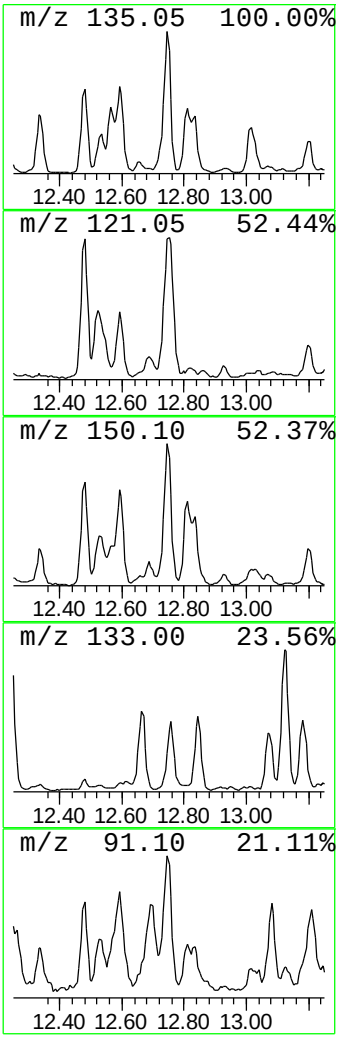
Peak Number 19 Phenol, 2,3,5,6-tetramethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.75	7.18 ng/ul	1478260	Acenaphthene-d10	14.33

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 2,3,5,6-tetramethyl-	150	C10H14O	000527-35-5	83
2	Phenol, 2,3,5,6-tetramethyl-	150	C10H14O	000527-35-5	83
3	Cyclohexanone, 2-(2-butynyl)-	150	C10H14O	054166-48-2	80
4	Phenol, 2-ethyl-4,5-dimethyl-	150	C10H14O	002219-78-5	76
5	Thymol	150	C10H14O	000089-83-8	74





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Sample : J5298-06
Misc :
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Instrument : BNA_M
ClientSampleId : E62T9

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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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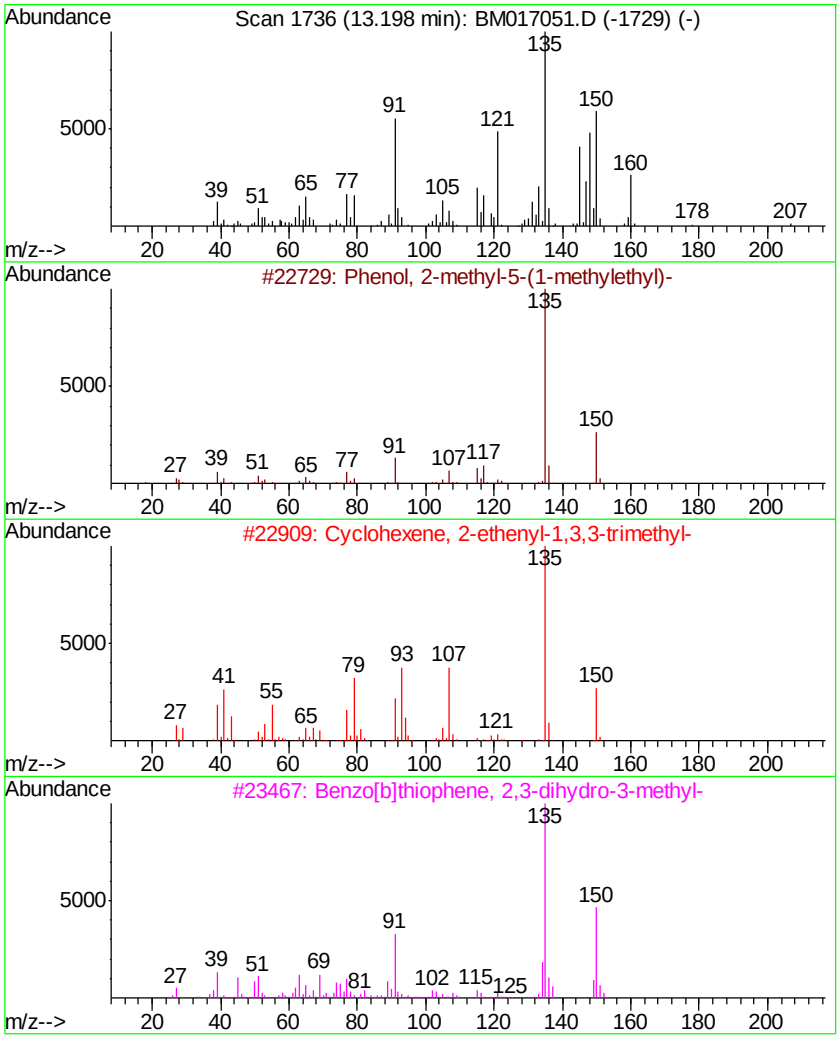
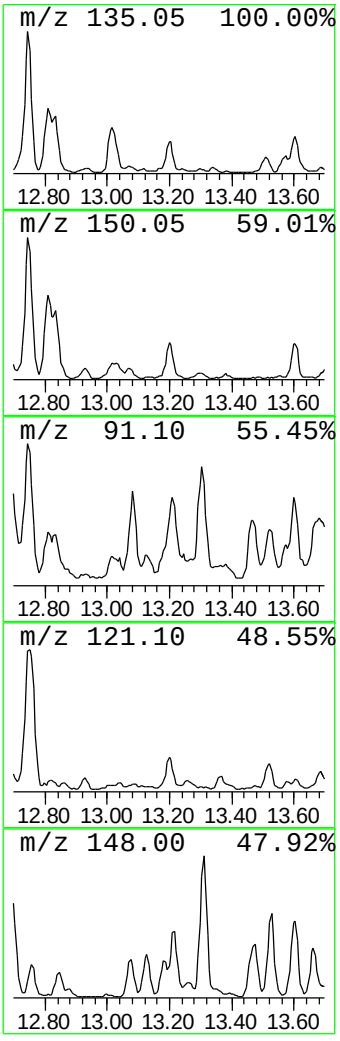
Peak Number 20 unknown-02 Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.20	2.61 ng/ul	536133	Acenaphthene-d10	14.33

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 2-methyl-5-(1-methylethyl)-	150	C10H14O	000499-75-2	38
2	Cyclohexene, 2-ethenyl-1,3,3-tri...	150	C11H18	005293-90-3	38
3	Benzo[b]thiophene, 2,3-dihydro-3...	150	C9H10S	006383-15-9	38
4	1-(4-Hydroxymethylphenyl)ethanone	150	C9H10O2	075633-63-5	38
5	Phenol, 2,3,4,6-tetramethyl-	150	C10H14O	003238-38-8	30





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Misc :
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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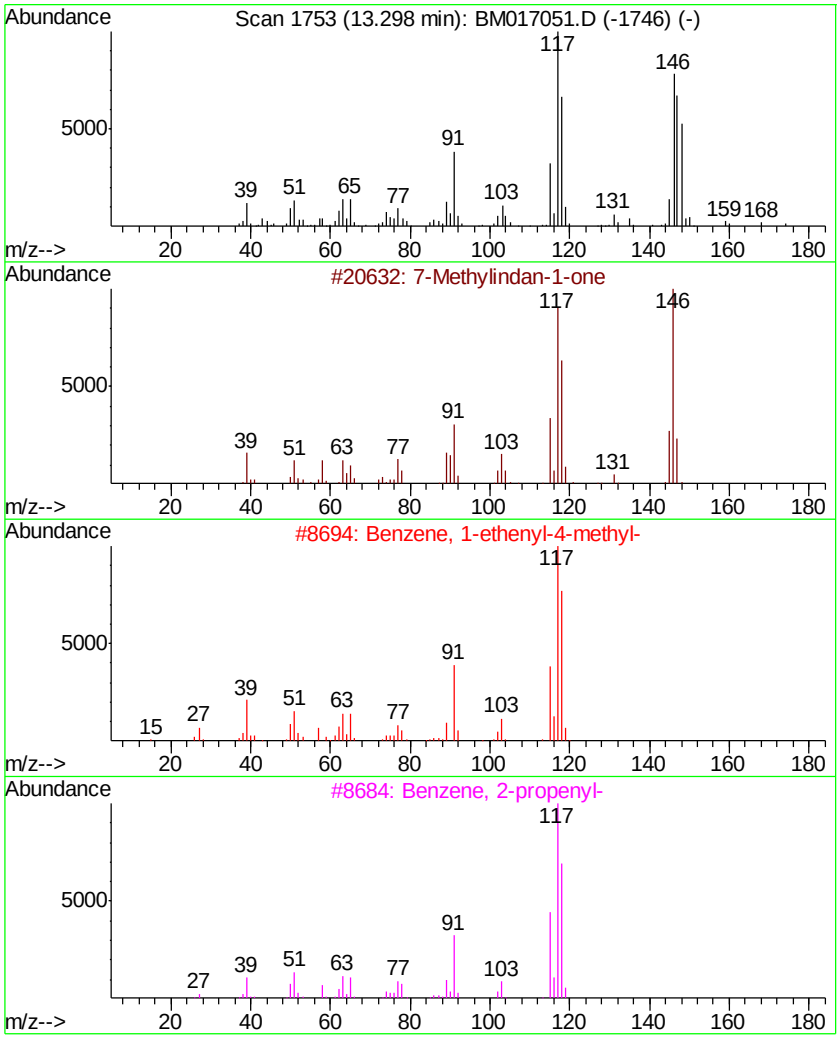
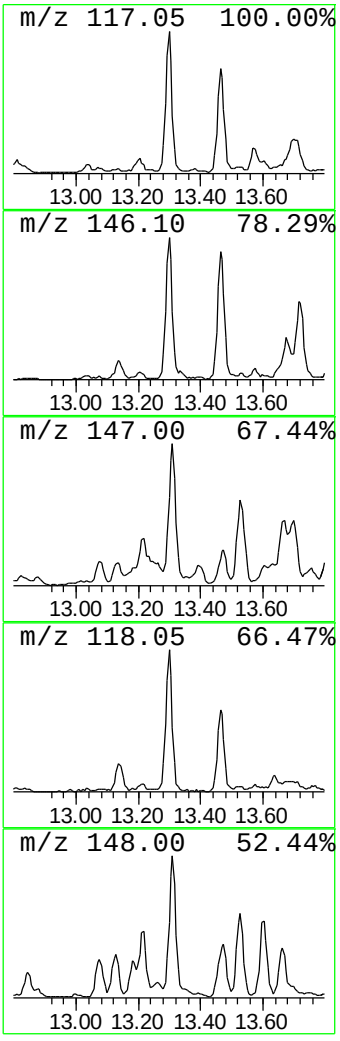
Peak Number 21 7-Methylindan-1-one Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.30	4.10 ng/ul	844112	Acenaphthene-d10	14.33

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	7-Methylindan-1-one	146	C10H10O	039627-61-7	70
2	Benzene, 1-ethenyl-4-methyl-	118	C9H10	000622-97-9	56
3	Benzene, 2-propenyl-	118	C9H10	000300-57-2	51
4	1,4-Benzenedicarboxaldehyde, 2-m...	148	C9H8O2	027587-17-3	46
5	Benzene, 2-propenyl-	118	C9H10	000300-57-2	42





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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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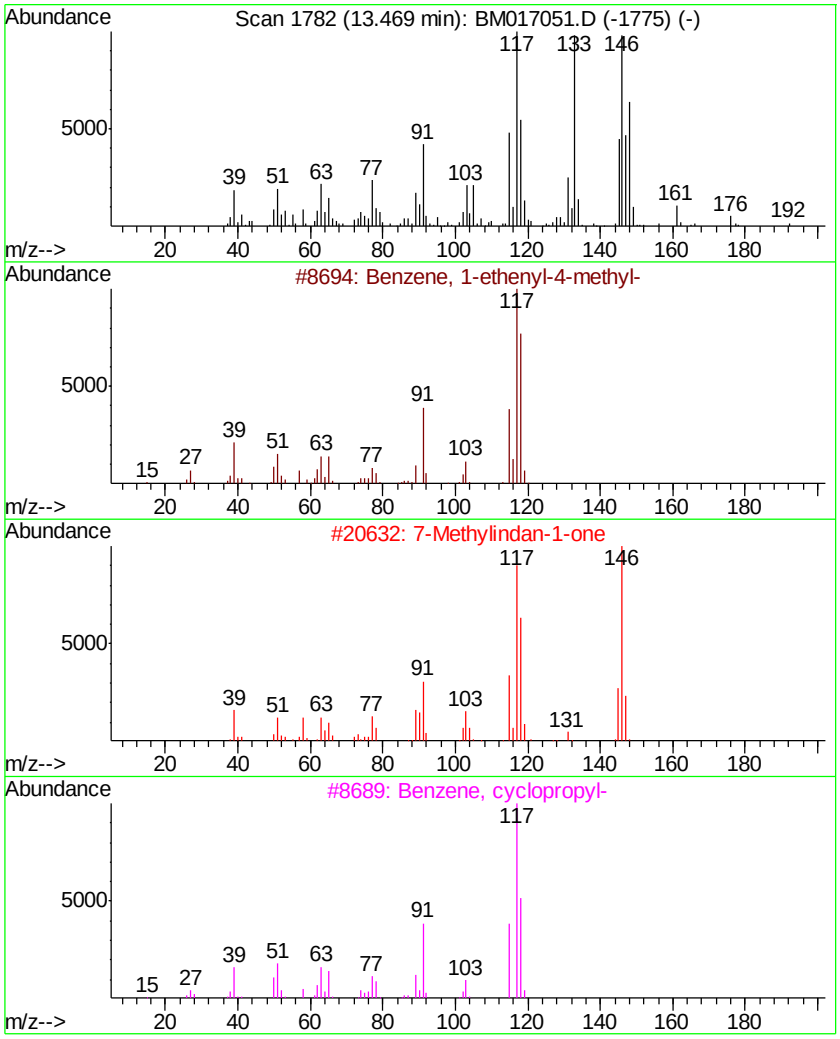
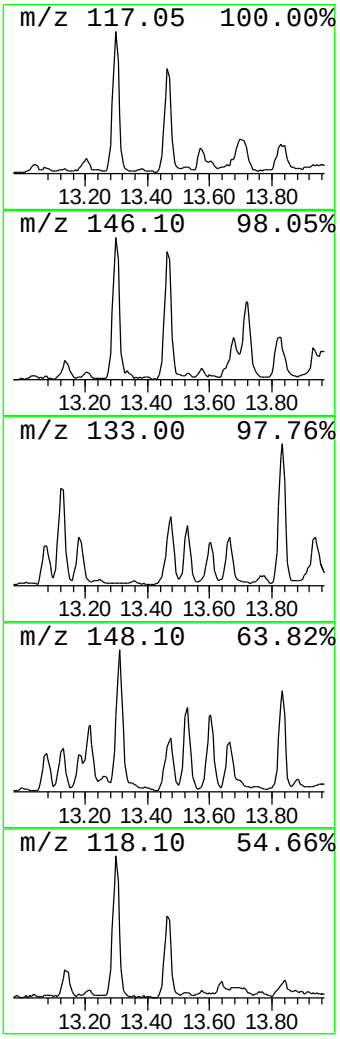
Peak Number 22 Benzene, 1-ethenyl-4-methyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.47	3.54 ng/ul	727544	Acenaphthene-d10	14.33

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethenyl-4-methyl-	118	C9H10	000622-97-9	55
2	7-Methylindan-1-one	146	C10H10O	039627-61-7	55
3	Benzene, cyclopropyl-	118	C9H10	000873-49-4	53
4	Naphth[1,2-b]oxirene, 1a,2,3,7b-...	146	C10H10O	002461-34-9	46
5	1H-Imidazole, 4,5-dihydro-2-phenyl-	146	C9H10N2	000936-49-2	38





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Data File : BM017051.D
Acq On : 06 Oct 2018 00:41
Operator : MJ/SJ
Sample : J5298-06
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : E62T9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM100418MA.M
Quant Title : SVOA CALIBRATION

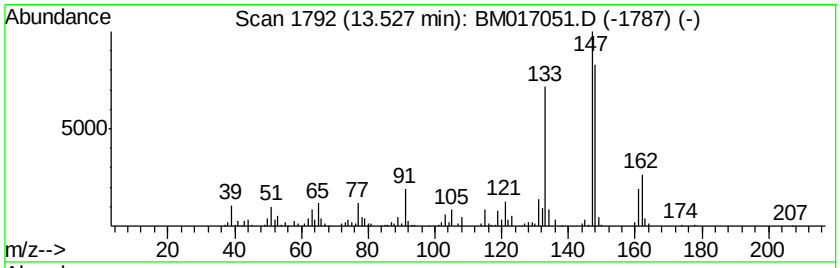
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

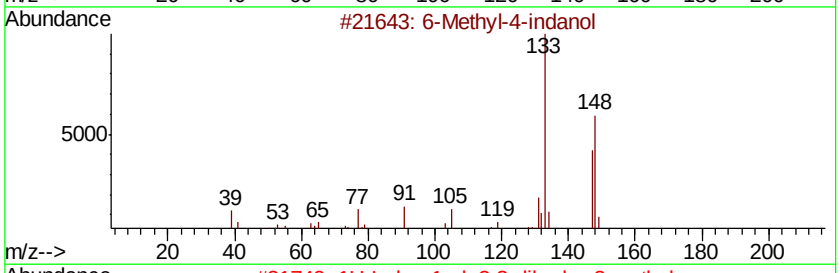
****

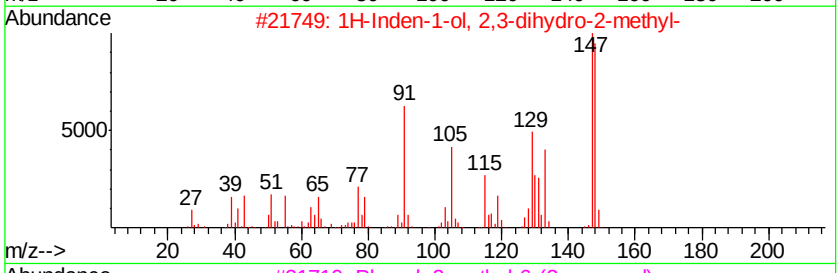
Peak Number 23 6-Methyl-4-indanol Concentration Rank 23

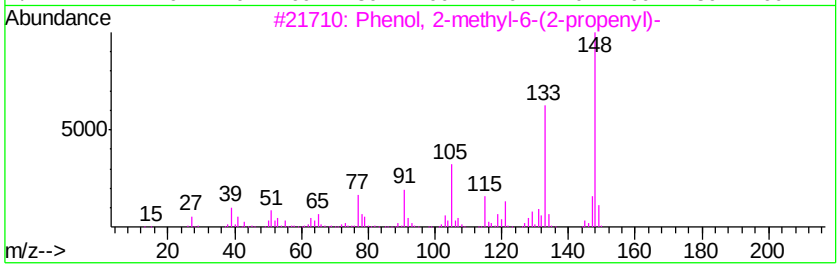
R.T.	EstConc	Area	Relative to ISTD	R.T.
13.53	2.57 ng/ul	529794	Acenaphthene-d10	14.33

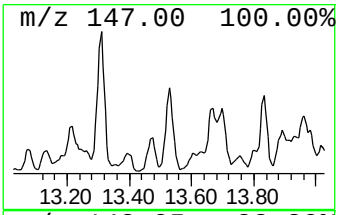
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	6-Methyl-4-indanol	148	C10H12O	020294-32-0	52
2	1H-Inden-1-ol, 2,3-dihydro-2-met...	148	C10H12O	017496-18-3	50
3	Phenol, 2-methyl-6-(2-propenyl)-	148	C10H12O	003354-58-3	49
4	Benzo[b]thiophene, 6-methyl-	148	C9H8S	016587-47-6	49
5	2-Allyl-4-methylphenol	148	C10H12O	006628-06-4	49

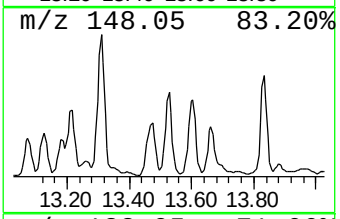


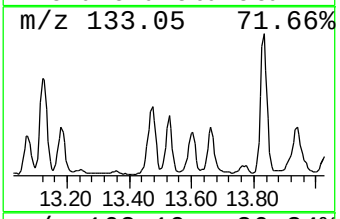


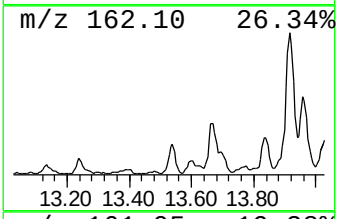


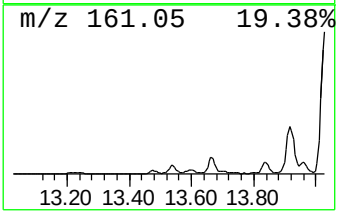












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Data File : BM017051.D
Acq On : 06 Oct 2018 00:41
Operator : MJ/SJ
Sample : J5298-06
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : E62T9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM100418MA.M
Quant Title : SVOA CALIBRATION

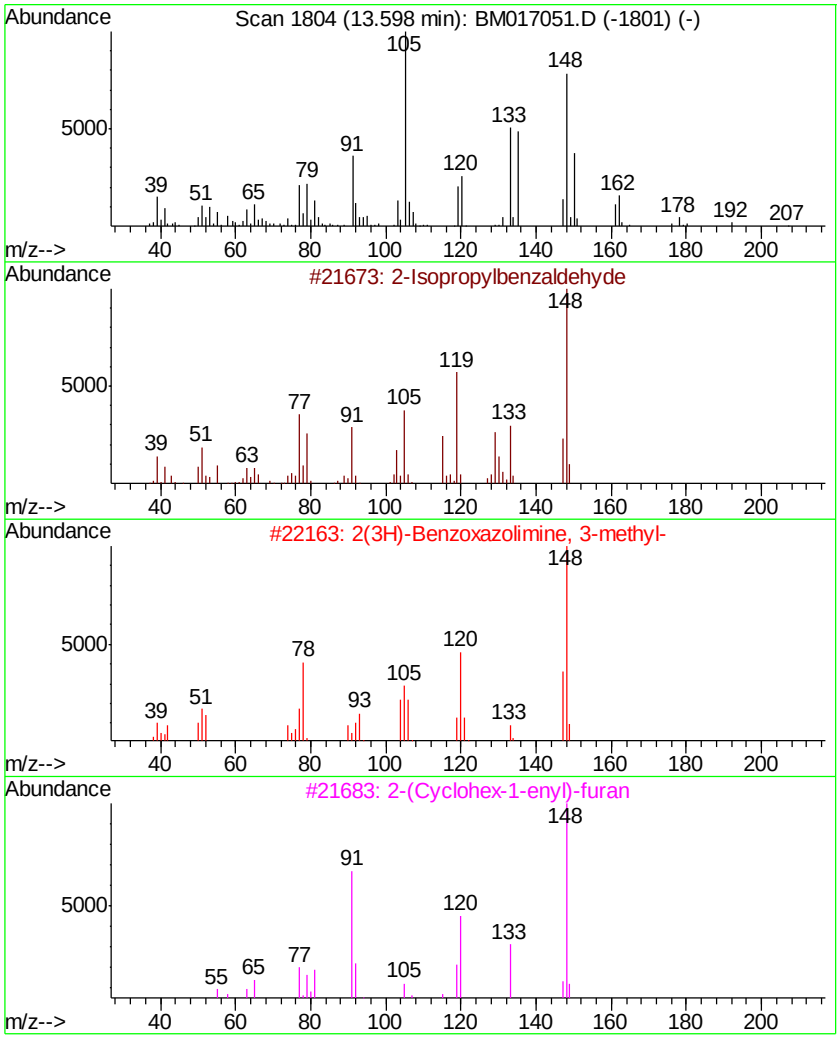
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

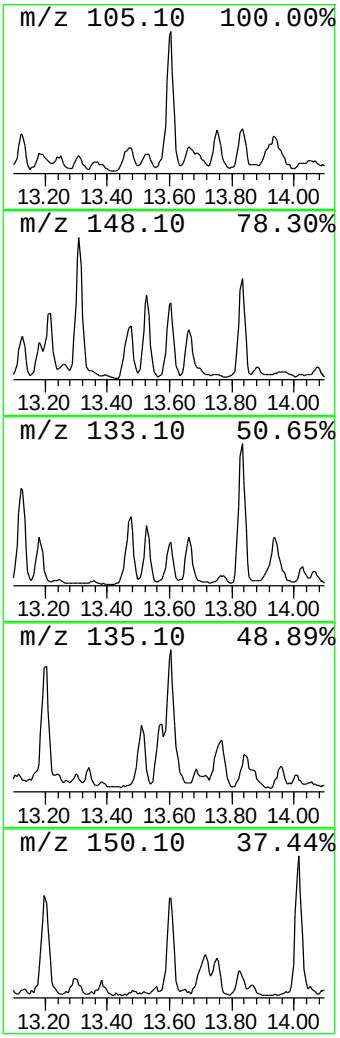
****

Peak Number 24 unknown-03 Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.60	3.07 ng/ul	631875	Acenaphthene-d10	14.33

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Isopropylbenzaldehyde	148	C10H12O	006502-22-3	41
2	2(3H)-Benzoxazolimine, 3-methyl-	148	C8H8N2O	018034-93-0	38
3	2-(Cyclohex-1-enyl)-furan	148	C10H12O	014174-50-6	38
4	3H-Indazol-3-one, 1,2-dihydro-1-...	148	C8H8N2O	001006-19-5	30
5	Pyrrole-3-carbonitrile, 5-formyl...	148	C8H8N2O	032487-71-1	25





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Data File : BM017051.D
Acq On : 06 Oct 2018 00:41
Operator : MJ/SJ
Sample : J5298-06
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : E62T9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM100418MA.M
Quant Title : SVOA CALIBRATION

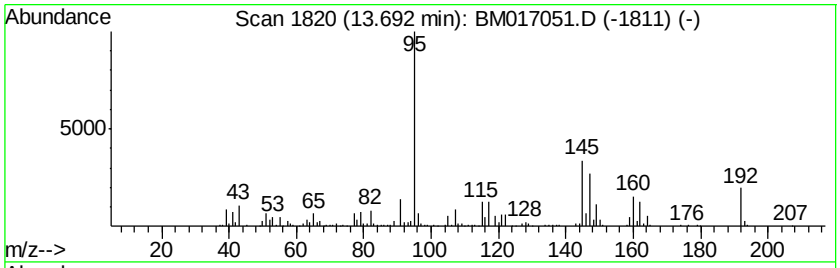
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

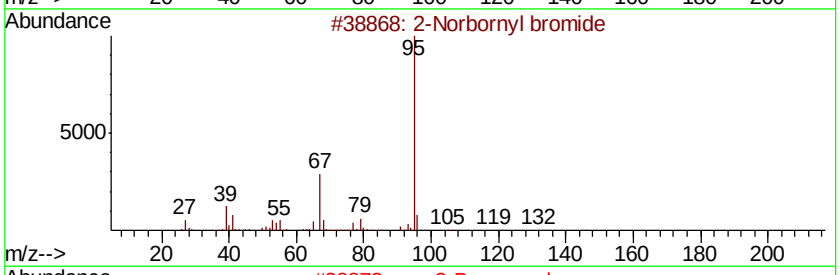
****

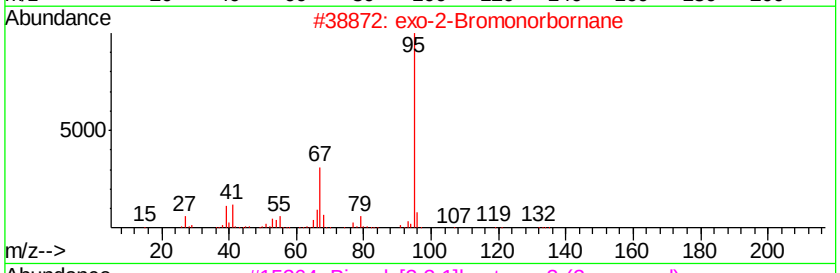
Peak Number 25 unknown-04 Concentration Rank 3

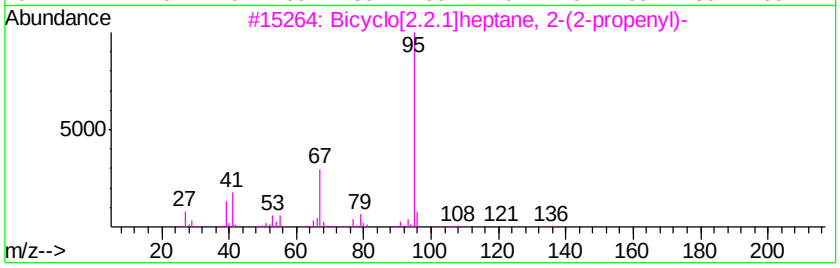
R.T.	EstConc	Area	Relative to ISTD	R.T.
13.69	7.95 ng/ul	1635610	Acenaphthene-d10	14.33

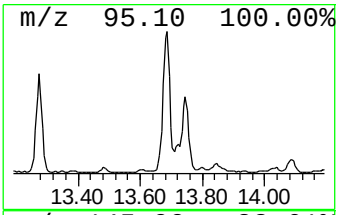
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Norbornyl bromide	174	C7H11Br	029342-65-2	38
2	exo-2-Bromonorbornane	174	C7H11Br	002534-77-2	38
3	Bicyclo[2.2.1]heptane, 2-(2-prop...	136	C10H16	002633-80-9	38
4	Borneol	154	C10H18O	000507-70-0	37
5	Furan, 2-ethyl-5-methyl-	110	C7H10O	001703-52-2	32

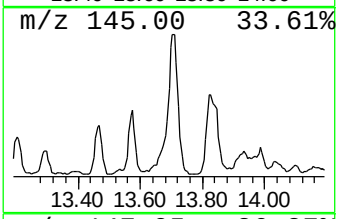


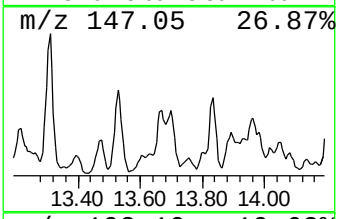


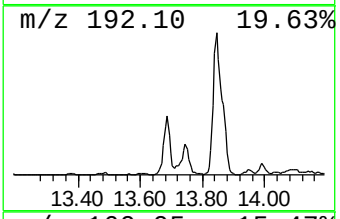


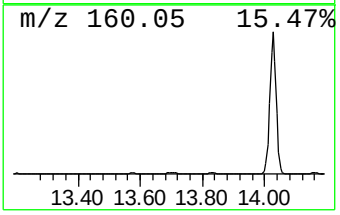












Library Search Compound Report

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Data File : BM017051.D
Acq On : 06 Oct 2018 00:41
Operator : MJ/SJ
Sample : J5298-06
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : E62T9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM100418MA.M
Quant Title : SVOA CALIBRATION

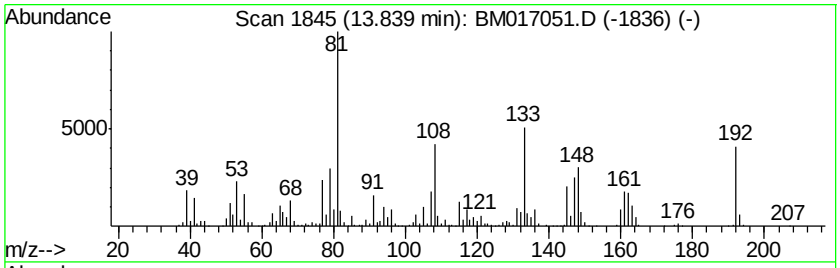
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

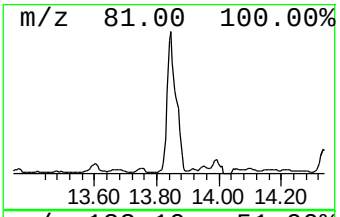
****

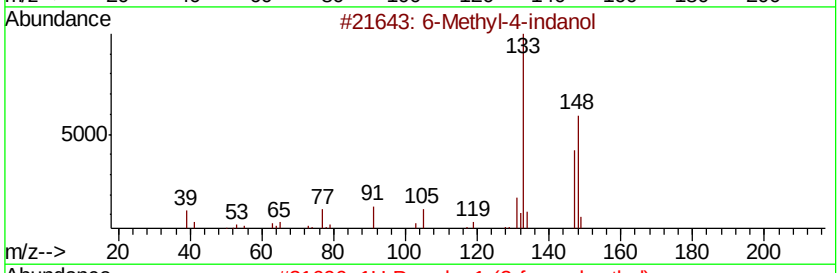
Peak Number 26 unknown-05 Concentration Rank 1

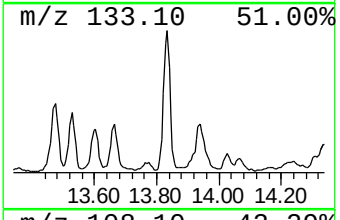
R.T.	EstConc	Area	Relative to ISTD	R.T.
13.84	10.16 ng/ul	2090210	Acenaphthene-d10	14.33

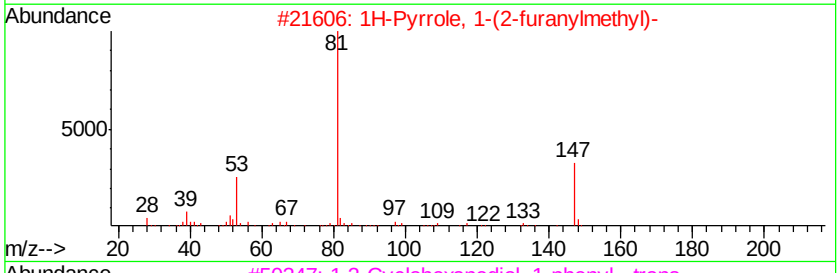
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	6-Methyl-4-indanol	148	C10H12O	020294-32-0	25
2	1H-Pyrrole, 1-(2-furanylmethyl)-	147	C9H9NO	001438-94-4	22
3	1,2-Cyclohexanediol, 1-phenyl-, ...	192	C12H16O2	027167-34-6	20
4	2-Pentyne, 4,4-dimethyl-	96	C7H12	000999-78-0	15
5	Cyclopropane, 1-chloro-2-ethenyl...	116	C6H9Cl	062337-93-3	11

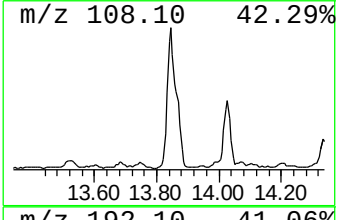


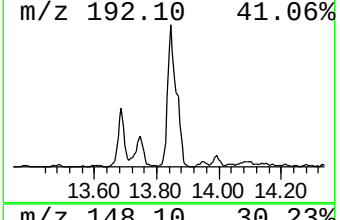


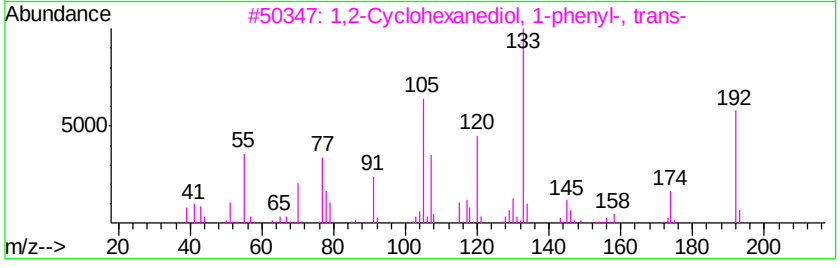


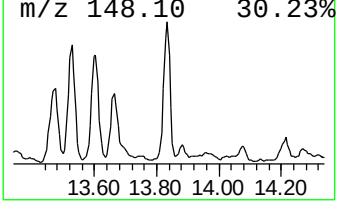












Library Search Compound Report

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Data File : BM017051.D
Acq On : 06 Oct 2018 00:41
Operator : MJ/SJ
Sample : J5298-06
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : E62T9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM100418MA.M
Quant Title : SVOA CALIBRATION

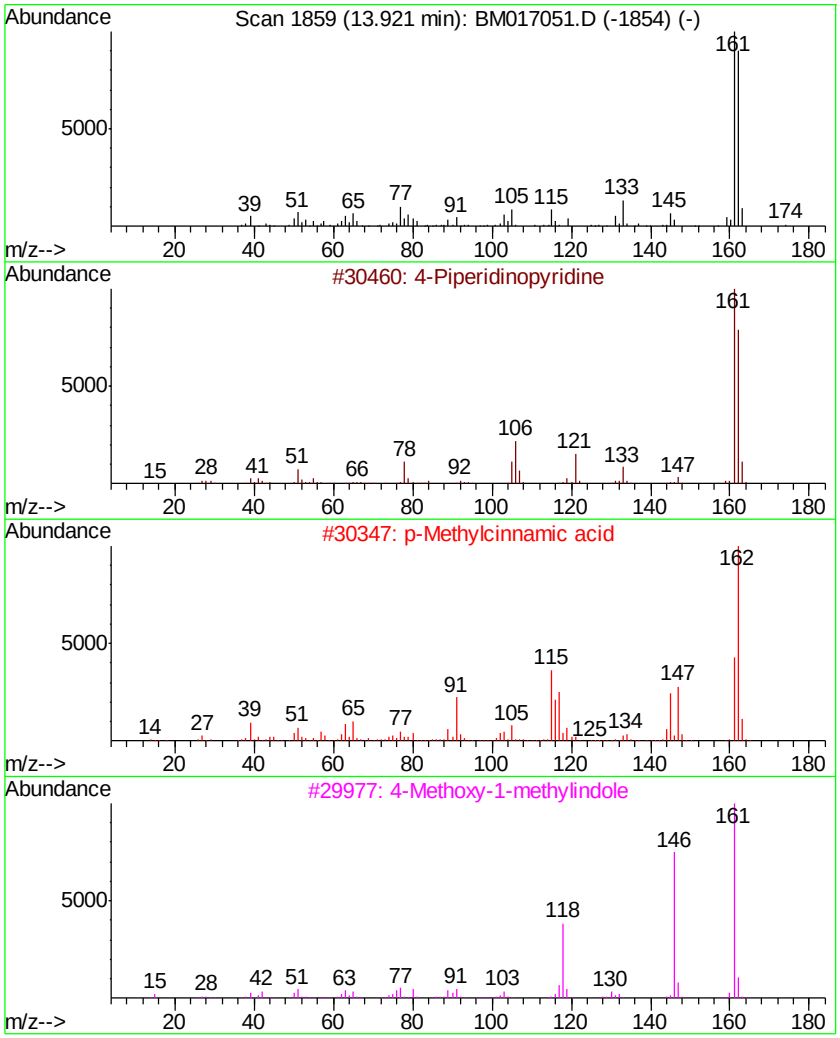
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

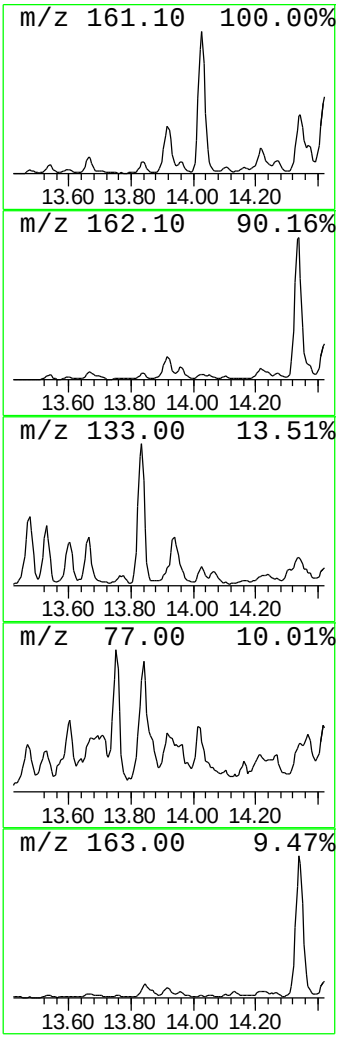
****

Peak Number 27 4-Piperidinopyridine Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.92	3.19 ng/ul	655902	Acenaphthene-d10	14.33

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-Piperidinopyridine	162	C10H14N2	002767-90-0	72
2	p-Methylcinnamic acid	162	C10H10O2	001866-39-3	43
3	4-Methoxy-1-methylindole	161	C10H11NO	007556-35-6	38
4	Furan, 2-(2-furanylmethyl)-5-met...	162	C10H10O2	013678-51-8	38
5	1H-Isoindole-1,3(2H)-dione, 2-am...	162	C8H6N2O2	001875-48-5	35





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Acq On : 06 Oct 2018 00:41
Operator : MJ/SJ
Sample : J5298-06
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : E62T9

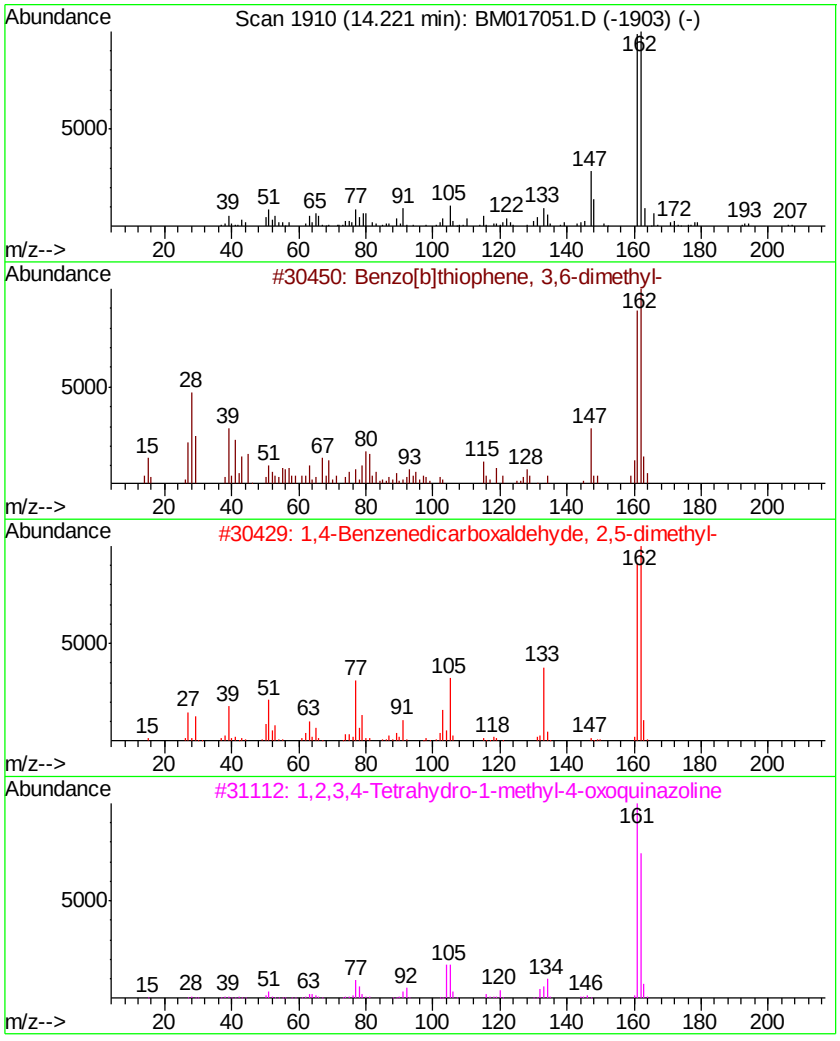
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM100418MA.M
Quant Title : SVOA CALIBRATION

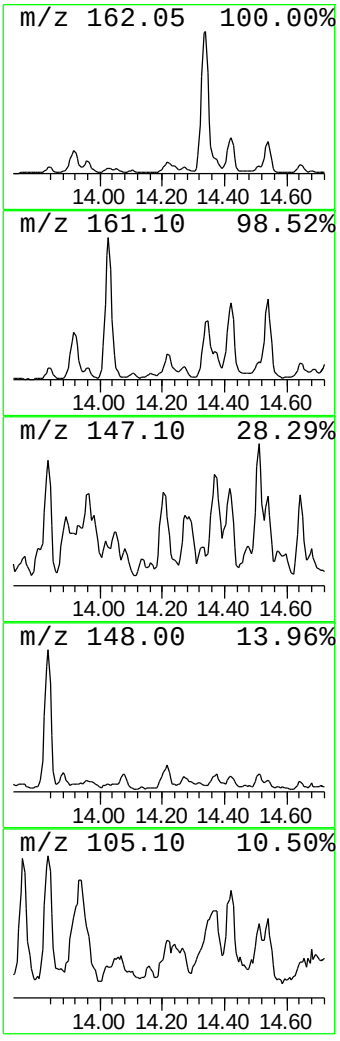
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 28 Benzo[b]thiophene, 3,6-dime... Concentration Rank 30

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.22	2.01 ng/ul	413971	Acenaphthene-d10	14.33

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzo[b]thiophene, 3,6-dimethyl-	162	C10H10S	016587-50-1	72
2	1,4-Benzenedicarboxaldehyde, 2,5...	162	C10H10O2	007044-92-0	53
3	1,2,3,4-Tetrahydro-1-methyl-4-ox...	162	C9H10N2O	035042-16-1	53
4	Benzo[b]thiophene, 2-ethyl-	162	C10H10S	001196-81-2	43
5	4-Butylbenzoic acid, 2-.+/-.chlo...	288	C17H17ClO2	1000293-59-7	38





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Acq On : 06 Oct 2018 00:41
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Sample : J5298-06
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : E62T9

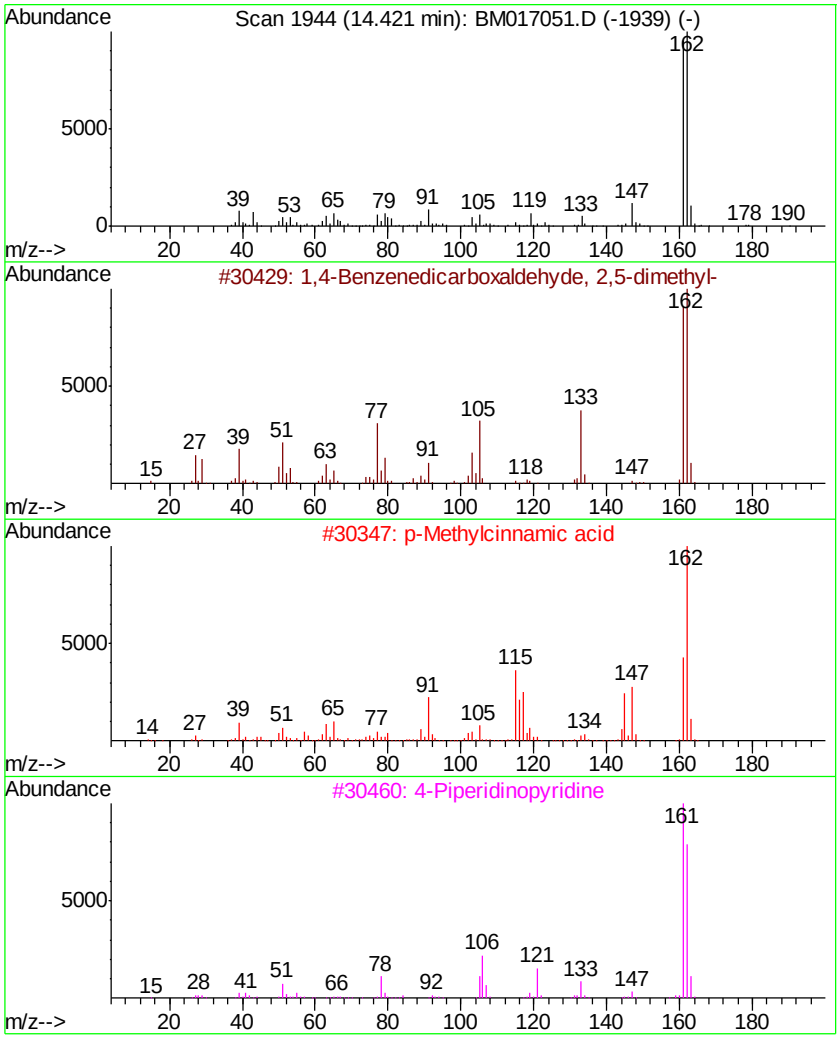
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM100418MA.M
Quant Title : SVOA CALIBRATION

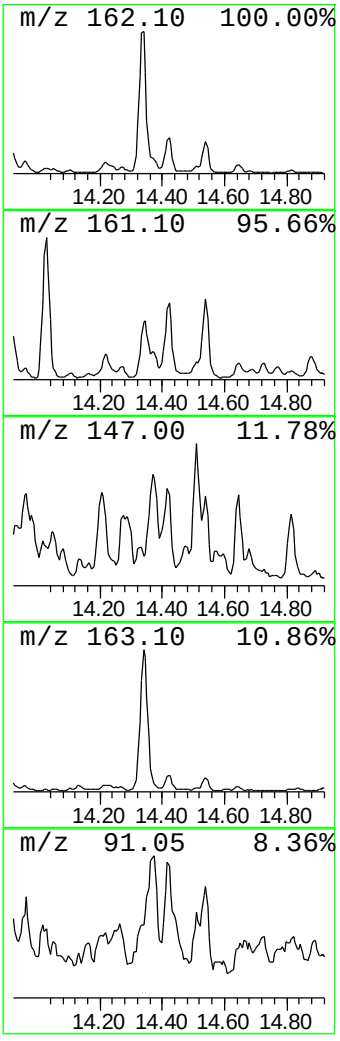
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 29 1,4-Benzenedicarboxaldehyde... Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.42	2.84 ng/ul	585171	Acenaphthene-d10	14.33

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,4-Benzenedicarboxaldehyde, 2,5...	162	C10H10O2	007044-92-0	64
2	p-Methylcinnamic acid	162	C10H10O2	001866-39-3	42
3	4-Piperidinopyridine	162	C10H14N2	002767-90-0	39
4	Boron, diethyl[3-imino-2-(1-imin...	191	C10H18BN3	136704-95-5	38
5	Isothiazole, 4-phenyl-	161	C9H7NS	000936-46-9	38





Library Search Compound Report

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Data File : BM017051.D
Acq On : 06 Oct 2018 00:41
Operator : MJ/SJ
Sample : J5298-06
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : E62T9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM100418MA.M
Quant Title : SVOA CALIBRATION

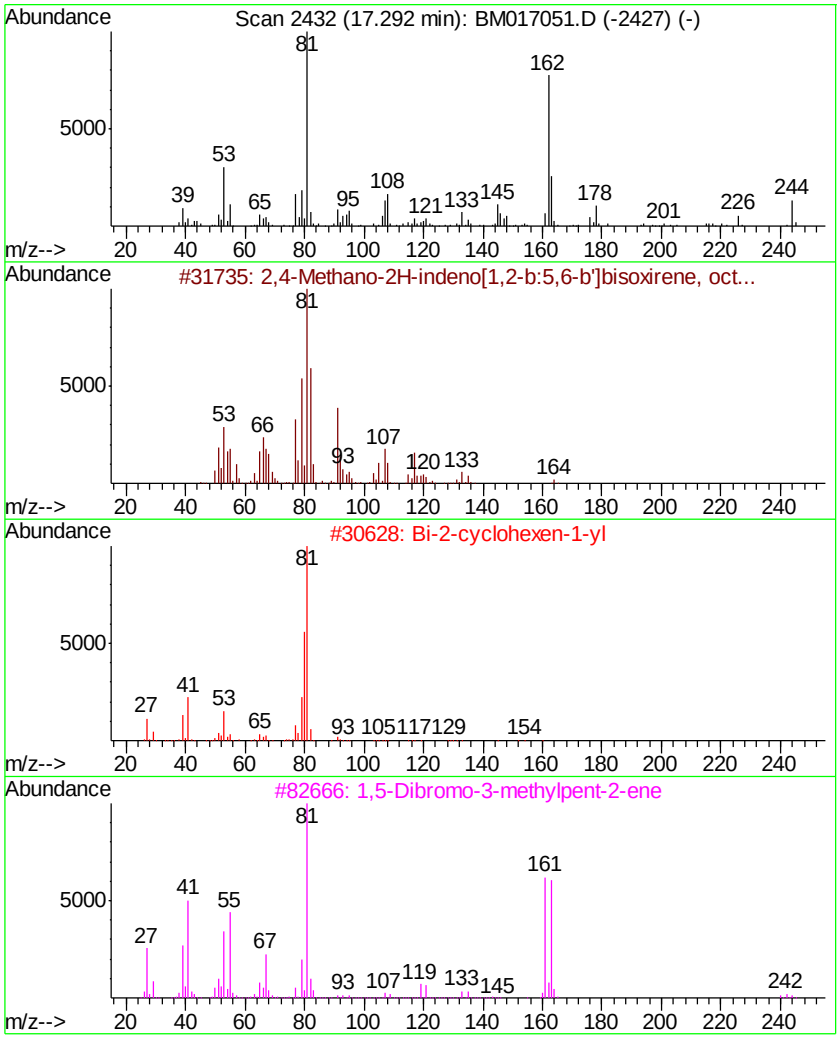
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

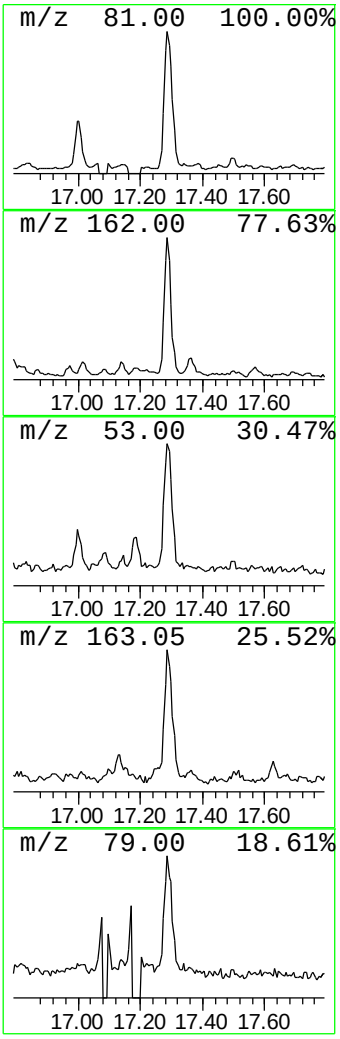
****

Peak Number 30 unknown-06 Concentration Rank 29

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.29	2.03 ng/ul	388096	Phenanthrene-d10	17.09

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,4-Methano-2H-indeno[1,2-b:5,6-...	164	C10H12O2	000081-21-0	43
2	Bi-2-cyclohexen-1-yl	162	C12H18	001541-20-4	38
3	1,5-Dibromo-3-methylpent-2-ene	240	C6H10Br2	133545-67-2	37
4	Cyclopropane, 1-methyl-1-isoprop...	96	C7H12	003422-07-9	35
5	1,E-4,Z-8-Dodecatriene	164	C12H20	083489-22-9	35





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100518\
Data File : BM017051.D
Acq On : 06 Oct 2018 00:41
Operator : MJ/SJ
Sample : J5298-06
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : E62T9

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM100418MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

					Internal Standard		
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Indane	8.06	2.9	ng/ul	227790	1	7.70	1557570	20.0
(DEL) Alkane: Cyc...	8.66	2.4	ng/ul	188417	1	7.70	1557570	20.0
Benzofuran, 2,3-d...	8.76	3.8	ng/ul	297477	1	7.70	1557570	20.0
Benzofuran, 2-met...	9.05	3.9	ng/ul	306228	1	7.70	1557570	20.0
Phenol, 2-ethyl-5...	10.30	2.5	ng/ul	342104	2	10.47	2732430	20.0
Phenol, 2,3,6-tri...	11.06	8.9	ng/ul	1221740	2	10.47	2732430	20.0
Phenol, 2-ethyl-6...	11.35	3.6	ng/ul	491057	2	10.47	2732430	20.0
Phenol, 2-(1-meth...	11.59	3.7	ng/ul	499900	2	10.47	2732430	20.0
Thymol	11.89	2.2	ng/ul	299989	2	10.47	2732430	20.0
Phenol, 3,4,5-tri...	12.19	6.7	ng/ul	920166	2	10.47	2732430	20.0
Phenol, 2-methyl-...	12.59	3.9	ng/ul	800242	3	14.33	4115980	20.0
unknown-01	12.69	6.5	ng/ul	1342330	3	14.33	4115980	20.0
Phenol, 2,3,5,6-t...	12.75	7.2	ng/ul	1478260	3	14.33	4115980	20.0
unknown-02	13.20	2.6	ng/ul	536133	3	14.33	4115980	20.0
7-Methylindan-1-one	13.30	4.1	ng/ul	844112	3	14.33	4115980	20.0
Benzene, 1-etheny...	13.47	3.5	ng/ul	727544	3	14.33	4115980	20.0
6-Methyl-4-indanol	13.53	2.6	ng/ul	529794	3	14.33	4115980	20.0
unknown-03	13.60	3.1	ng/ul	631875	3	14.33	4115980	20.0
unknown-04	13.69	8.0	ng/ul	1635610	3	14.33	4115980	20.0
unknown-05	13.84	10.2	ng/ul	2090210	3	14.33	4115980	20.0
4-Piperidinopyridine	13.92	3.2	ng/ul	655902	3	14.33	4115980	20.0
Benzo[b]thiophene...	14.22	2.0	ng/ul	413971	3	14.33	4115980	20.0
1,4-Benzenedicarb...	14.42	2.8	ng/ul	585171	3	14.33	4115980	20.0
unknown-06	17.29	2.0	ng/ul	388096	4	17.09	3832330	20.0