LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG061917\
Data File : BG027547.D
Acq On : 20 Jun 2017 2:56
Operator : SJ/MA
Sample : I3627-13 25X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : F5E24

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

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 1

Method : Z:\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG061617.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	8.518	778	788	799	rVB2	127642	233477	3.22%	0.426%
2	8.894	841	852	860	rVB	94438	172871	2.38%	0.315%
3	9.059	873	880	888	rBV	79864	147882	2.04%	0.270%
4	10.722	1155	1163	1175	rBV3	50203	121713	1.68%	0.222%
5	11.338	1260	1268	1271	rVV	198233	396337	5.46%	0.723%
6	11.397	1271	1278	1291	rVV	3592125	7259311	100.00%	13.238%
7	11.509	1291	1297	1305	rVB2	119714	231821	3.19%	0.423%
8	12.954	1531	1543	1554	rVB	1164137	2050946	28.25%	3.740%
9	13.172	1570	1580	1588	rVV	556544	956511	13.18%	1.744%
10	13.935	1705	1710	1722	rVV	277964	460200	6.34%	0.839%
11	14.135	1731	1744	1755	rVB	112560	240019	3.31%	0.438%
12	14.276	1755	1768	1777	rBV3	205952	566615	7.81%	1.033%
13	14.423	1786	1793	1798	rBV	272611	536436	7.39%	0.978%
14	14.476	1798	1802	1809	rVV2	196661	332211	4.58%	0.606%
15	14.658	1825	1833	1845	rBV	127623	265975	3.66%	0.485%
16	14.823	1850	1861	1868	rBV4	91845	193942	2.67%	0.354%
17	15.058	1894	1901	1905	rBV	158680	271489	3.74%	0.495%
18	15.105	1905	1909	1914	rVV2	329156	547400	7.54%	0.998%
19	15.169	1914	1920	1926	rVV	1298489	2117836	29.17%	3.862%
20	15.228	1926	1930	1936	rVV	49512	80423	1.11%	0.147%
21	15.299	1936	1942	1951	rVV	50423	118716	1.64%	0.216%
22	15.410	1951	1961	1968	rVV2	63543	154097	2.12%	0.281%
23	15.498	1968	1976	1982	rVV	916335	1557792	21.46%	2.841%
24	15.563	1982	1987	1996	rVB	77382	126497	1.74%	0.231%
25	15.704	2004	2011	2016	rBV2	60621	109908	1.51%	0.200%
26	15.757	2016	2020	2032	rVB	70255	114987	1.58%	0.210%
27	15.880	2032	2041	2043	rBV	55441	118154	1.63%	0.215%
28	16.051	2066	2070	2074	rBV2	45743	78065	1.08%	0.142%
29	16.150	2079	2087	2095	rVB	1362710	2213039	30.49%	4.036%
30	16.239	2095	2102	2104	rVV2	69528	113587	1.56%	0.207%
31	16.268	2104	2107	2110	rVV2	98176	171570	2.36%	0.313%
32	16.309	2110	2114	2119	rVV3	181951	325824	4.49%	0.594%
33	16.356	2119	2122	2125	rVV4	50262	78416	1.08%	0.143%
34	16.397	2125	2129	2131	rVV2	98173	153795	2.12%	0.280%

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

Instrument : BNA_G
ClientSampleId : F5E24

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

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 1

Method : Z:\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG061617.M
Title : SVOA CALIBRATION

35	16.432	2131	2135	2143	rVB	213414	377311	5.20%	0.688%
36	16.579	2143	2160	2169	rBV	233682	576325	7.94%	1.051%
37	16.679	2169	2177	2181	rVV2	47706	100320	1.38%	0.183%
38	16.785	2188	2195	2208	rVB4	80131	187918	2.59%	0.343%
39	16.938	2216	2221	2227	rVV	128181	192680	2.65%	0.351%
40	17.143	2244	2256	2261	rBV2	183878	459913	6.34%	0.839%
41	17.196	2261	2265	2271	rVV3	79571	142019	1.96%	0.259%
42	17.302	2276	2283	2287	rVB4	69739	143643	1.98%	0.262%
43	17.373	2287	2295	2298	rBV3	79167	183358	2.53%	0.334%
44	17.414	2298	2302	2310	rVB3	78281	157321	2.17%	0.287%
45	17.496	2311	2316	2322	rVV6	42176	109845	1.51%	0.200%
46	17.572	2322	2329	2340	rVV4	100372	360796	4.97%	0.658%
47	17.672	2340	2346	2361	rVB	304989	549530	7.57%	1.002%
48	17.854	2369	2377	2380	rBV	413259	714738	9.85%	1.303%
49	17.907	2380	2386	2393	rVV	3781735	6467763	89.10%	11.794%
50	17.989	2393	2400	2413	rVV	619899	1098279	15.13%	2.003%
51	18.254	2439	2445	2460	rBV	295859	609781	8.40%	1.112%
52	18.436	2471	2476	2487	rVB5	41190	99118	1.37%	0.181%
53	18.571	2494	2499	2509	rVB3	40557	91379	1.26%	0.167%
54	18.724	2519	2525	2529	rVV	314677	508305	7.00%	0.927%
55	18.771	2529	2533	2539	rVB	427839	641567	8.84%	1.170%
56	18.853	2539	2547	2551	rBV	172769	296010	4.08%	0.540%
57	18.924	2551	2559	2571	rVV2	481167	1147634	15.81%	2.093%
58	19.206	2602	2607	2614	rBV	214886	315784	4.35%	0.576%
59	19.447	2638	2648	2652	rBV5	61958	130901	1.80%	0.239%
60	19.499	2652	2657	2660	rVV	62157	98132	1.35%	0.179%
61	19.617	2668	2677	2682	rVV2	117502	255621	3.52%	0.466%
62	19.687	2682	2689	2697	rVV4	74886	233680	3.22%	0.426%
63	19.758	2697	2701	2710	rVB3	38028	92325	1.27%	0.168%
64	19.893	2716	2724	2730	rBV	2179857	3482627	47.97%	6.351%
65	20.134	2759	2765	2773	rVB2	71085	145745	2.01%	0.266%
66	20.216	2773	2779	2781	rBV2	420258	684205	9.43%	1.248%
67	20.252	2781	2785	2792	rVV	1394557	2258193	31.11%	4.118%
68	20.310	2792	2795	2802	rVB	99295	158832	2.19%	0.290%
69	20.422	2802	2814	2819	rBV	102299	178636	2.46%	0.326%
70	20.563	2829	2838	2844	rBV3	116505	295837	4.08%	0.539%
71	20.733	2853	2867	2876	rVB	381509	747243	10.29%	1.363%

LSC Area Percent Report

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Sample : I3627-13 25X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : F5E24

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

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Peak separation: 1

Method : Z:\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG061617.M
Title : SVOA CALIBRATION

72	20.833	2876	2884	2889	rBV	384416	638865	8.80%	1.165%
73	20.898	2889	2895	2899	rVV3	99801	198607	2.74%	0.362%
74	21.051	2915	2921	2924	rBV	56969	94121	1.30%	0.172%
75	21.098	2925	2929	2941	rVB	65832	147358	2.03%	0.269%
76	21.291	2957	2962	2969	rBV4	55461	110407	1.52%	0.201%
77	21.403	2978	2981	2992	rVB5	50051	115634	1.59%	0.211%
78	21.503	2992	2998	3004	rBV2	53140	127059	1.75%	0.232%
79	21.762	3032	3042	3047	rBV	113122	248040	3.42%	0.452%
80	21.803	3047	3049	3055	rVB	72371	111783	1.54%	0.204%
81	21.867	3055	3060	3064	rBV2	52257	90602	1.25%	0.165%
82	22.073	3086	3095	3100	rBV	35959	84202	1.16%	0.154%
83	22.214	3108	3119	3126	rBV3	641745	1956147	26.95%	3.567%
84	22.279	3126	3130	3144	rVB	349758	710248	9.78%	1.295%
85	22.455	3153	3160	3168	rBV2	80866	174719	2.41%	0.319%
86	22.684	3192	3199	3207	rBV3	63968	140681	1.94%	0.257%
87	23.042	3250	3260	3268	rVV	59752	176242	2.43%	0.321%
88	23.119	3269	3273	3282	rVB2	38320	80947	1.12%	0.148%
89	23.354	3308	3313	3317	rVV3	35094	79402	1.09%	0.145%
90	23.413	3317	3323	3331	rVB6	47242	123000	1.69%	0.224%
91	23.512	3332	3340	3348	rVB6	28886	76982	1.06%	0.140%
92	23.847	3388	3397	3414	rVV6	20786	104340	1.44%	0.190%
93	24.711	3531	3544	3553	rBV	147479	635221	8.75%	1.158%
94	25.005	3585	3594	3606	rVB	30865	105926	1.46%	0.193%
95	25.539	3674	3685	3696	rBV2	63238	210958	2.91%	0.385%
96	25.704	3703	3713	3725	rVB2	102428	343731	4.74%	0.627%
97	25.880	3730	3743	3753	rBV2	211729	751067	10.35%	1.370%
98	28.706	4205	4224	4237	rVB10	15817	76776	1.06%	0.140%
99	30.093	4445	4460	4477	rVB4	21526	129767	1.79%	0.237%
100	30.592	4532	4545	4563	rBV4	14103	84244	1.16%	0.154%

Sum of corrected areas: 54838252

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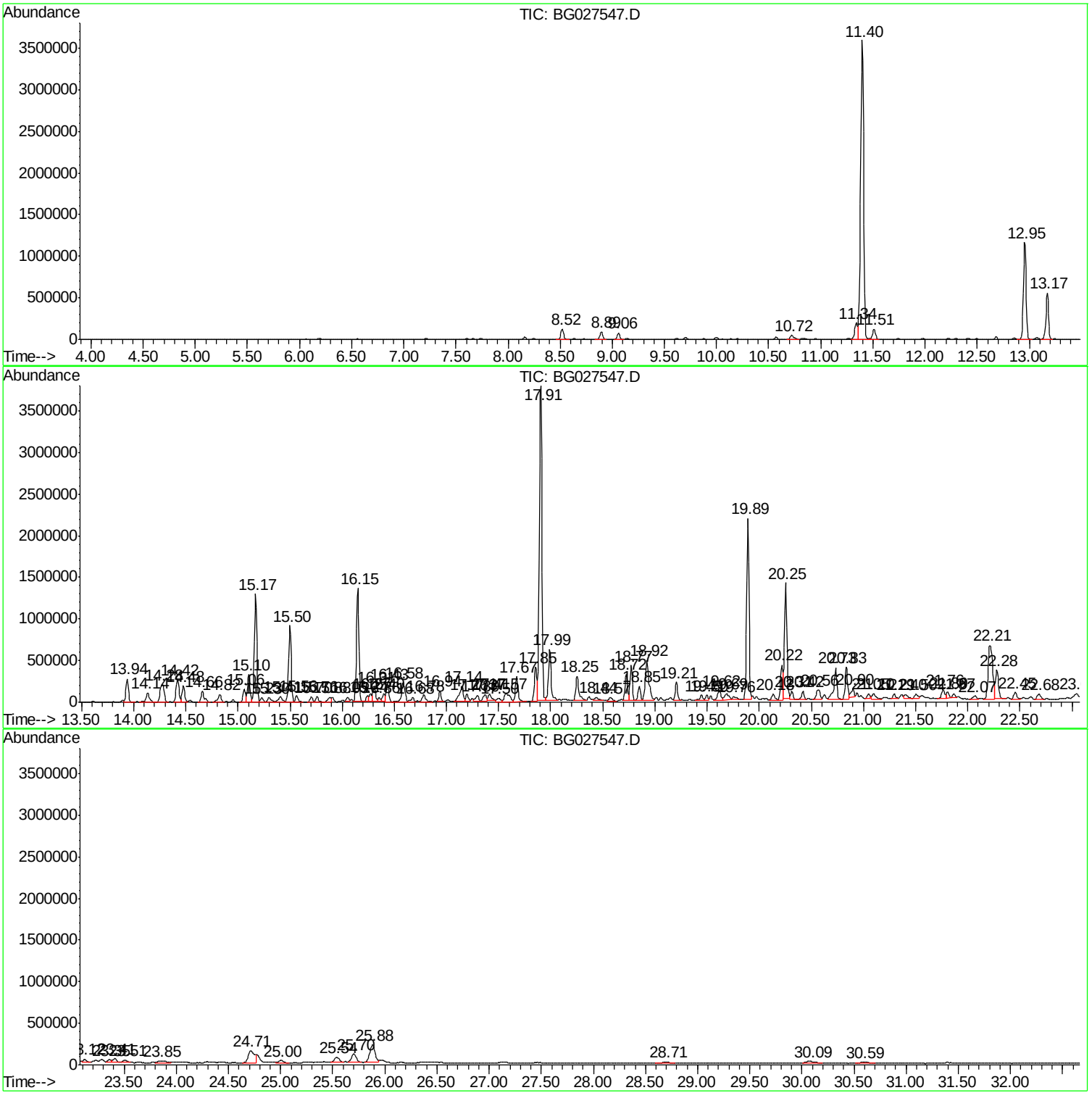
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : F5E24

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG061617.M
Quant Title : SVOA CALIBRATION

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



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Quant Title : SVOA CALIBRATION

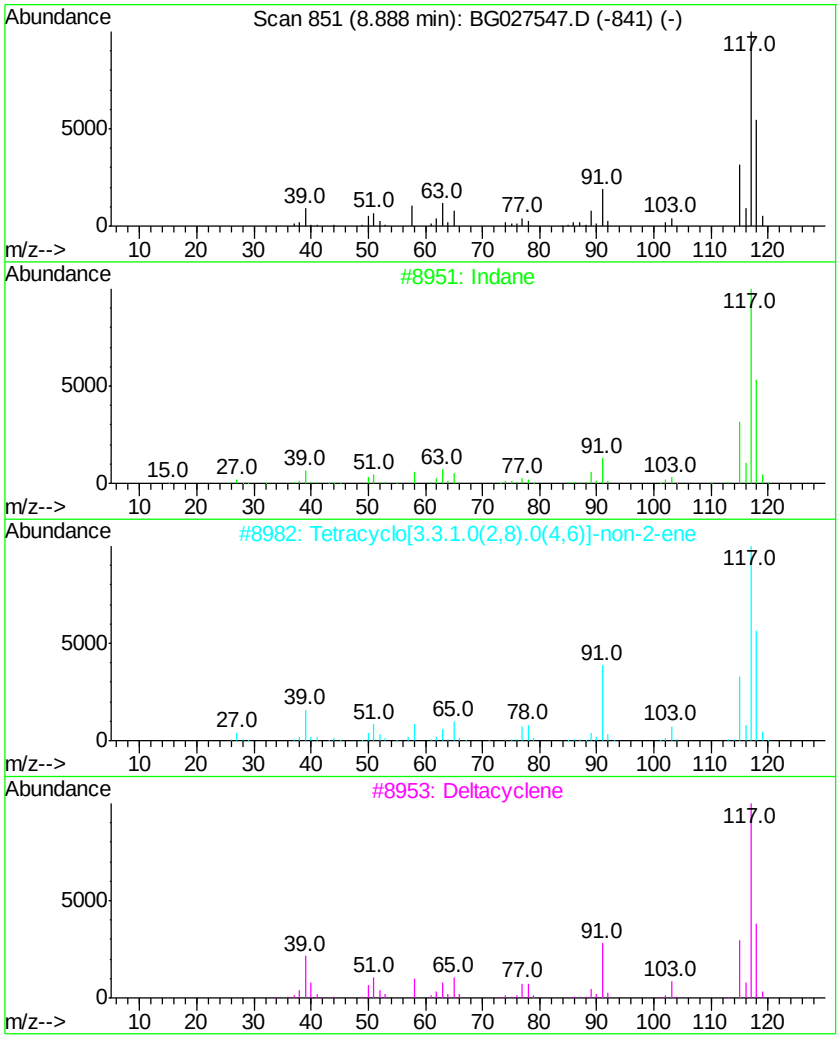
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

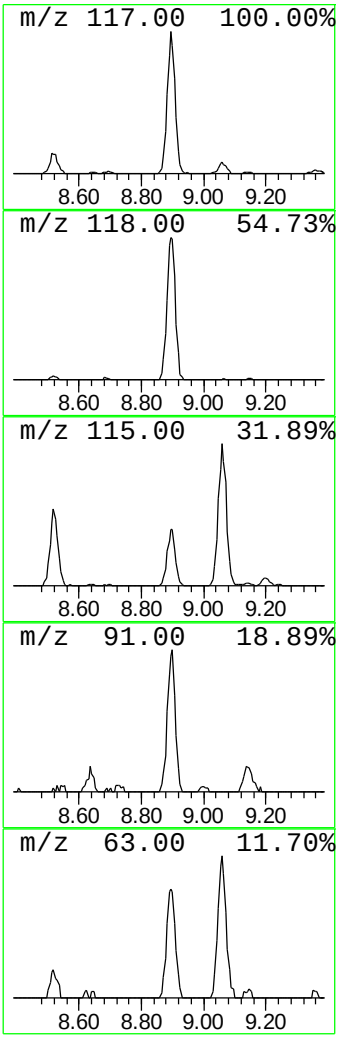
****

Peak Number 1 Indane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.89	14.81 ng/ul	172871	1,4-Dichlorobenzene-d4	8.52

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	91
2	Tetracyclo[3.3.1.0(2,8).0(4,6)]-...	118	C9H10	1000191-13-7	74
3	Deltacyclene	118	C9H10	007785-10-6	68
4	Indane	118	C9H10	000496-11-7	64
5	Indane	118	C9H10	000496-11-7	60





Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG061917\
Data File : BG027547.D
Acq On : 20 Jun 2017 2:56
Operator : SJ/MA
Sample : I3627-13 25X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : F5E24

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG061617.M
Quant Title : SVOA CALIBRATION

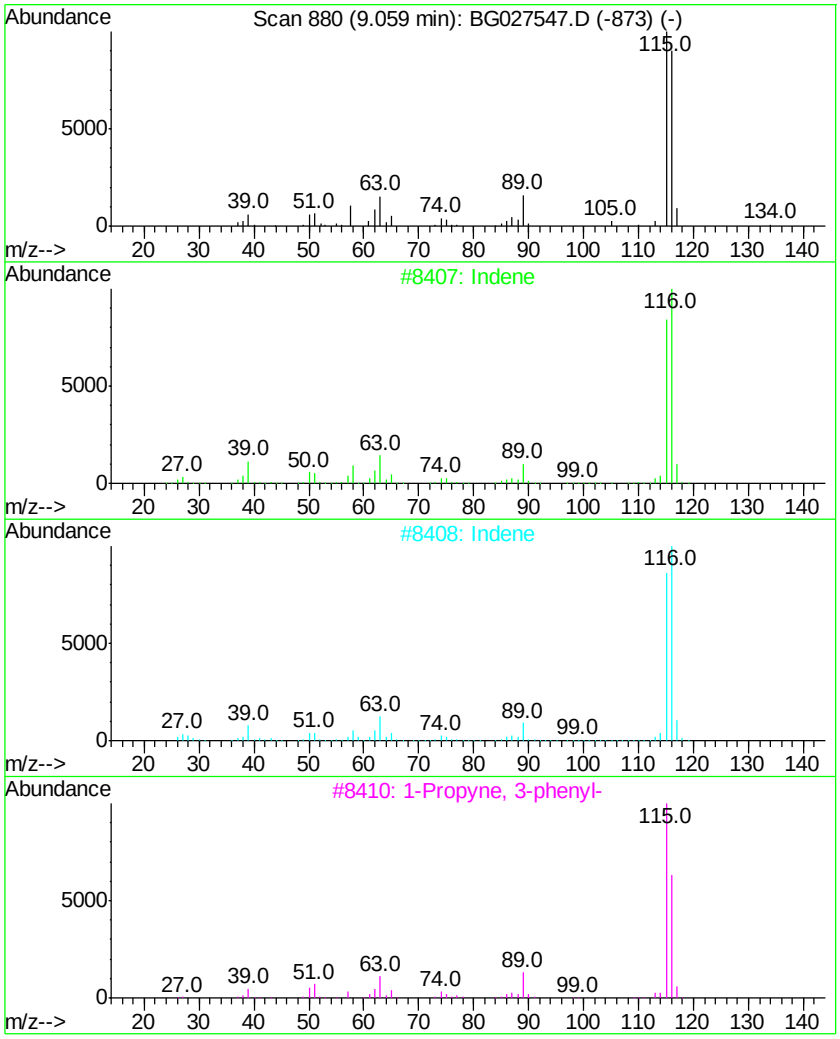
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

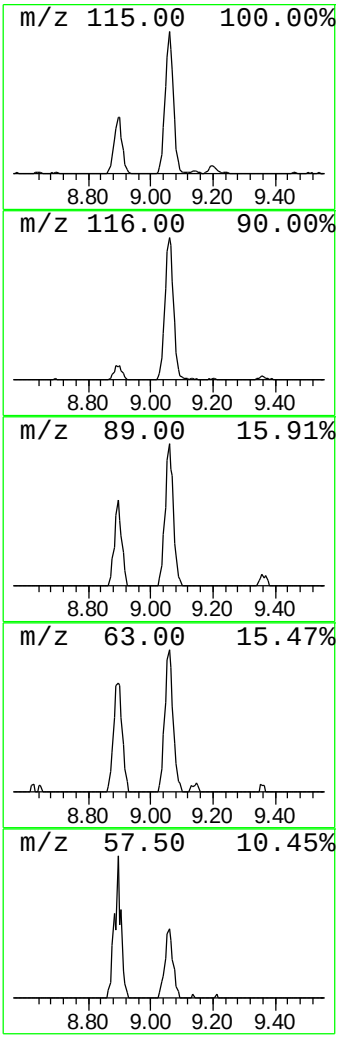
****

Peak Number 2 Indene Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.06	12.67 ng/ul	147882	1,4-Dichlorobenzene-d4	8.52

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indene	116	C9H8	000095-13-6	94
2	Indene	116	C9H8	000095-13-6	94
3	1-Propyne, 3-phenyl-	116	C9H8	010147-11-2	94
4	Benzene, 1-propynyl-	116	C9H8	000673-32-5	91
5	Benzene, 1-ethynyl-4-methyl-	116	C9H8	000766-97-2	91





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Instrument : BNA_G
ClientSampleId : F5E24

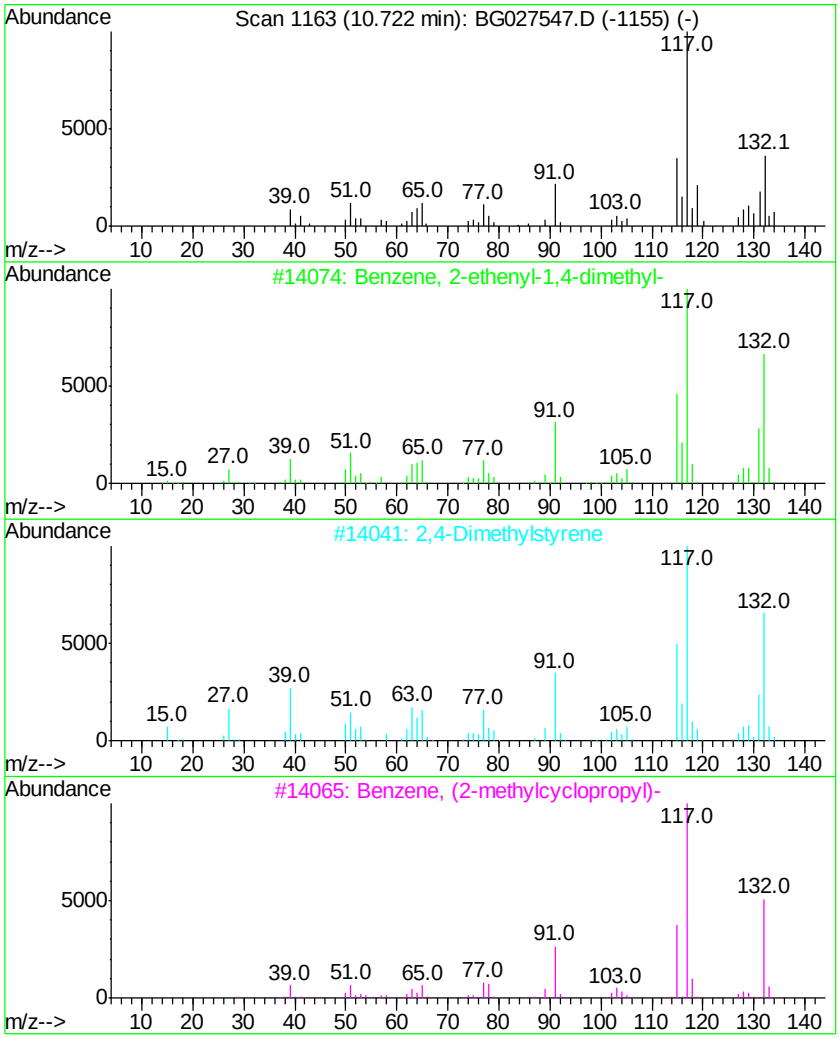
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG061617.M
Quant Title : SVOA CALIBRATION

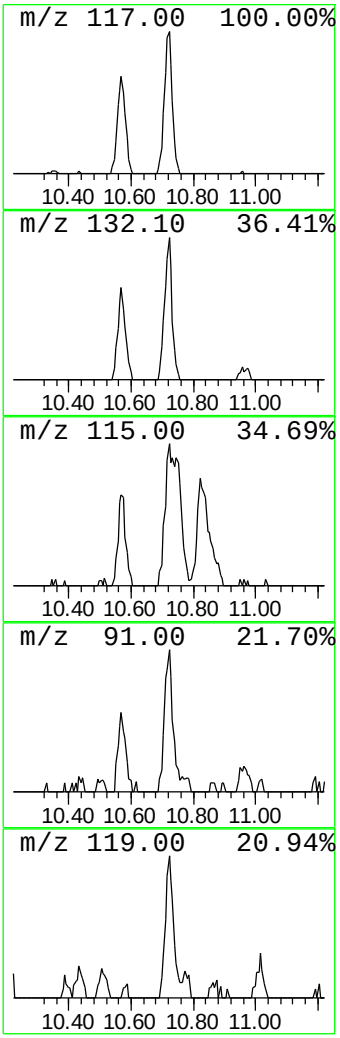
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.72	6.14 ng/ul	121713	Naphthalene-d8	11.34

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	93
2	2,4-Dimethylstyrene	132	C10H12	002234-20-0	93
3	Benzene, (2-methylcyclopropyl)-	132	C10H12	1000327-39-0	91
4	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	87
5	Benzene, 2-ethenyl-1,3-dimethyl-	132	C10H12	002039-90-9	87





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Sample : I3627-13 25X
Misc :
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Instrument : BNA_G
ClientSampleId : F5E24

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG061617.M
Quant Title : SVOA CALIBRATION

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

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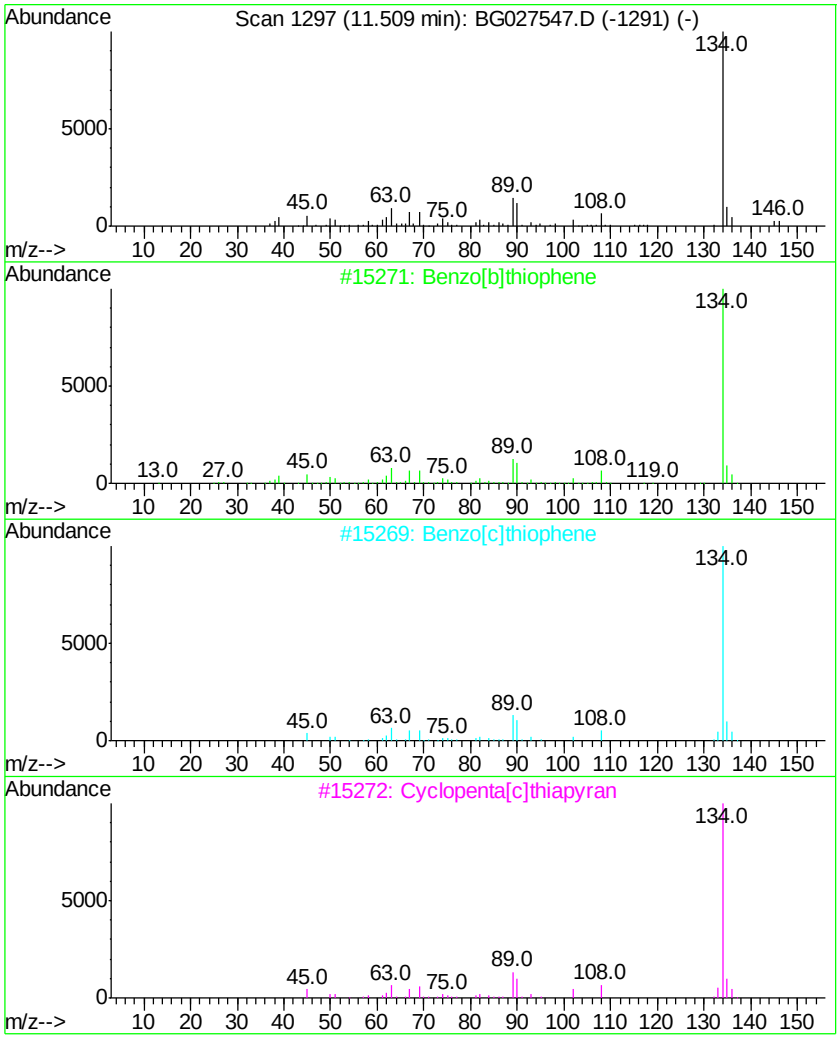
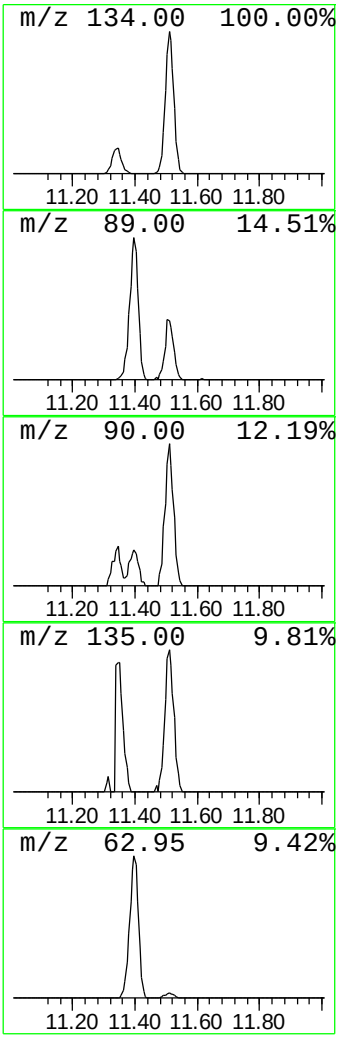
Peak Number 4 Benzo[b]thiophene Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.51	11.70 ng/ul	231821	Naphthalene-d8	11.34

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzo[b]thiophene	134	C8H6S	000095-15-8	97
2	Benzo[c]thiophene	134	C8H6S	000270-82-6	94
3	Cyclopenta[c]thiapyran	134	C8H6S	000270-63-3	91
4	Benzo[b]thiophene	134	C8H6S	000095-15-8	91
5	Benzo[b]thiophene	134	C8H6S	000095-15-8	91





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Data File : BG027547.D
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Operator : SJ/MA
Sample : I3627-13 25X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : F5E24

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG061617.M
Quant Title : SVOA CALIBRATION

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

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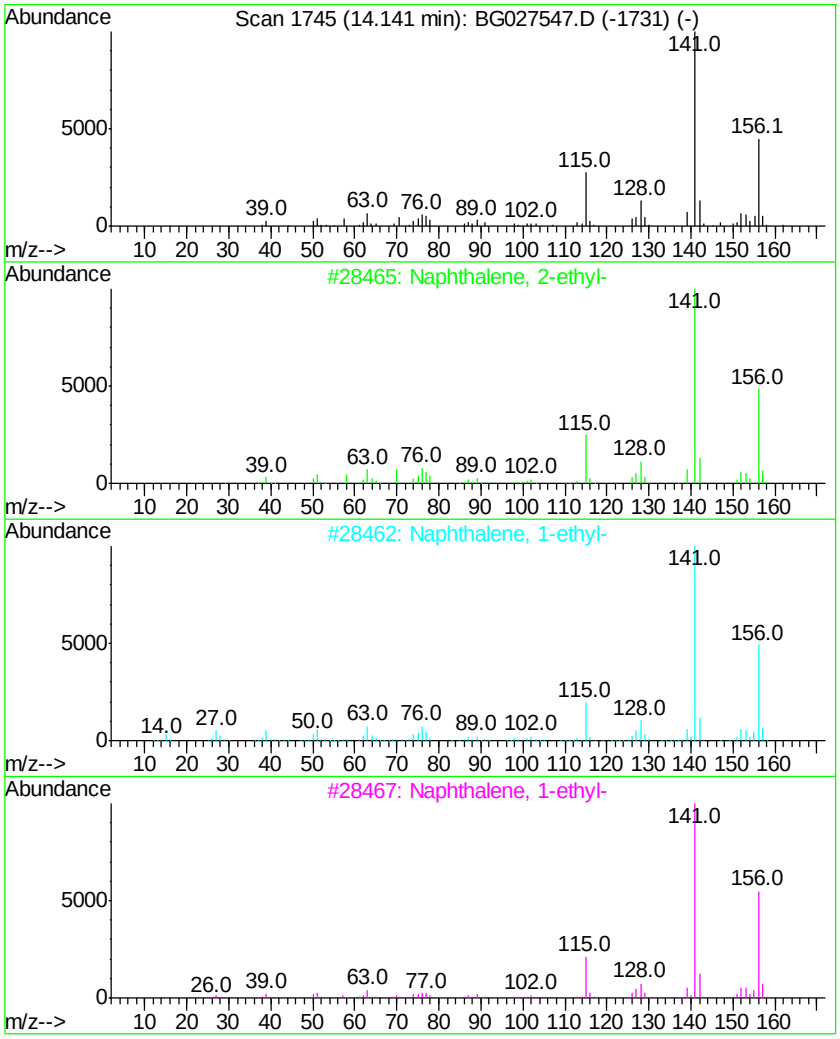
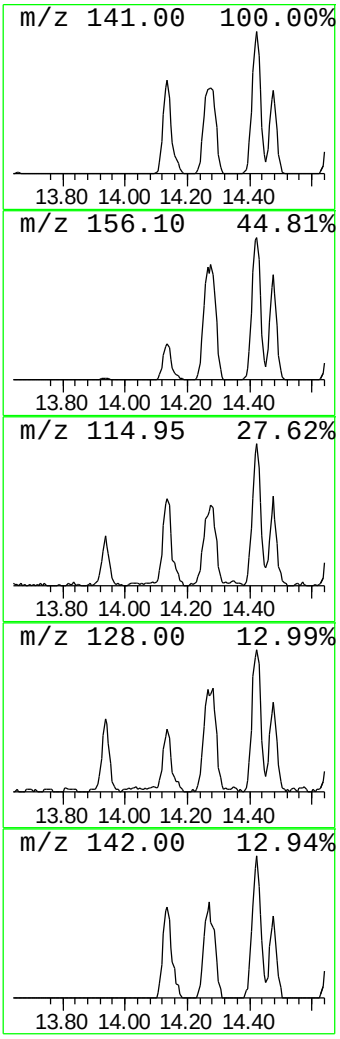
Peak Number 6 Naphthalene, 2-ethyl- Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.14	8.77 ng/ul	240019	Acenaphthene-d10	15.10

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2-ethyl-	156	C12H12	000939-27-5	96
2	Naphthalene, 1-ethyl-	156	C12H12	001127-76-0	94
3	Naphthalene, 1-ethyl-	156	C12H12	001127-76-0	94
4	Naphthalene, 2-ethyl-	156	C12H12	000939-27-5	91
5	Naphthalene, 2-ethyl-	156	C12H12	000939-27-5	91





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Data File : BG027547.D
Acq On : 20 Jun 2017 2:56
Operator : SJ/MA
Sample : I3627-13 25X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : F5E24

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG061617.M
Quant Title : SVOA CALIBRATION

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

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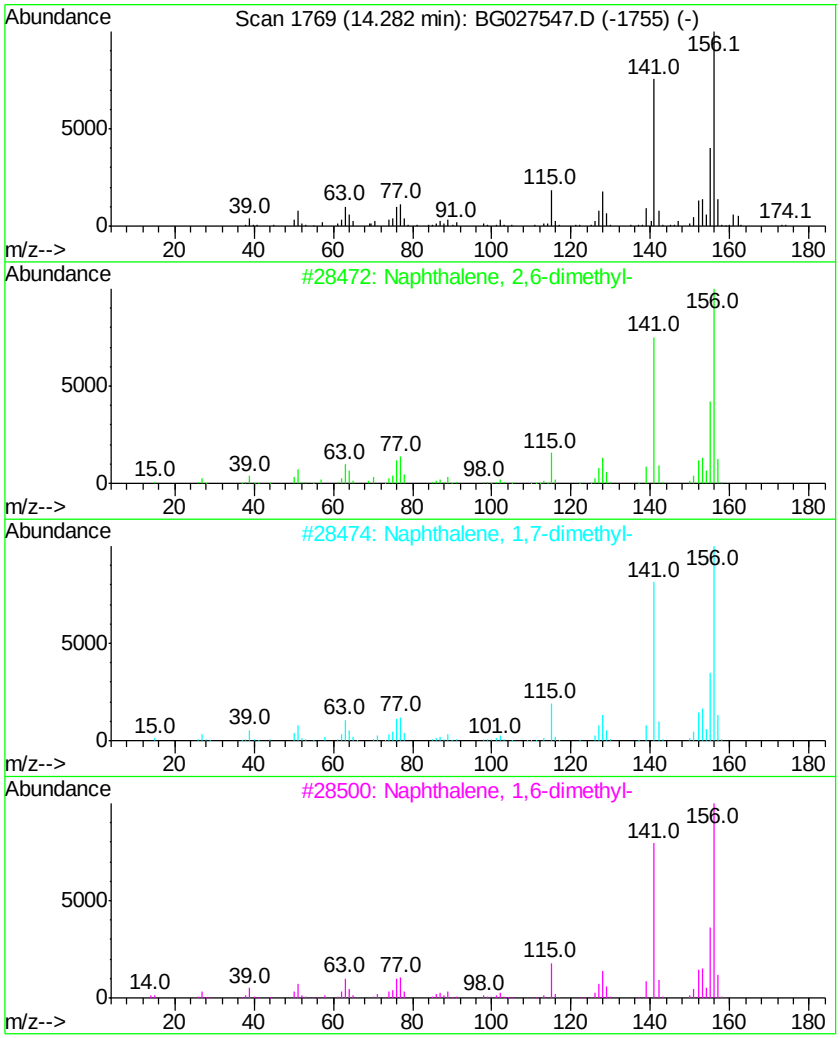
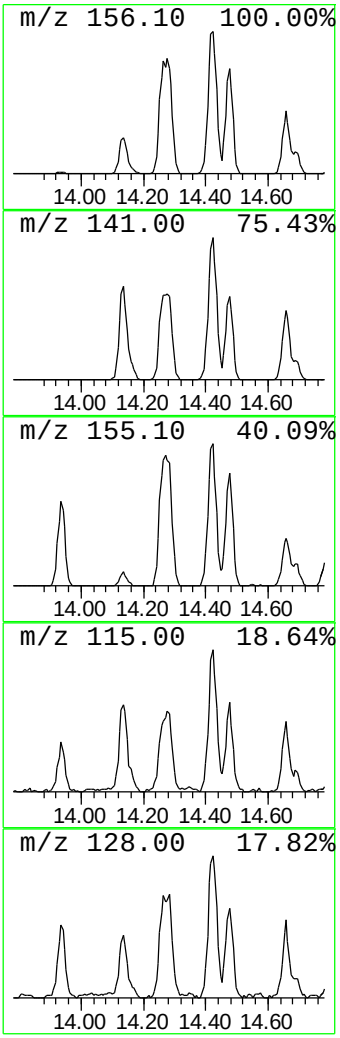
Peak Number 7 Naphthalene, 2,6-dimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.28	20.70 ng/ul	566615	Acenaphthene-d10	15.10

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	98
2	Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	98
3	Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	98
4	Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	96
5	Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	96





Library Search Compound Report

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Sample : I3627-13 25X
Misc :
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Instrument : BNA_G
ClientSampleId : F5E24

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Quant Title : SVOA CALIBRATION

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

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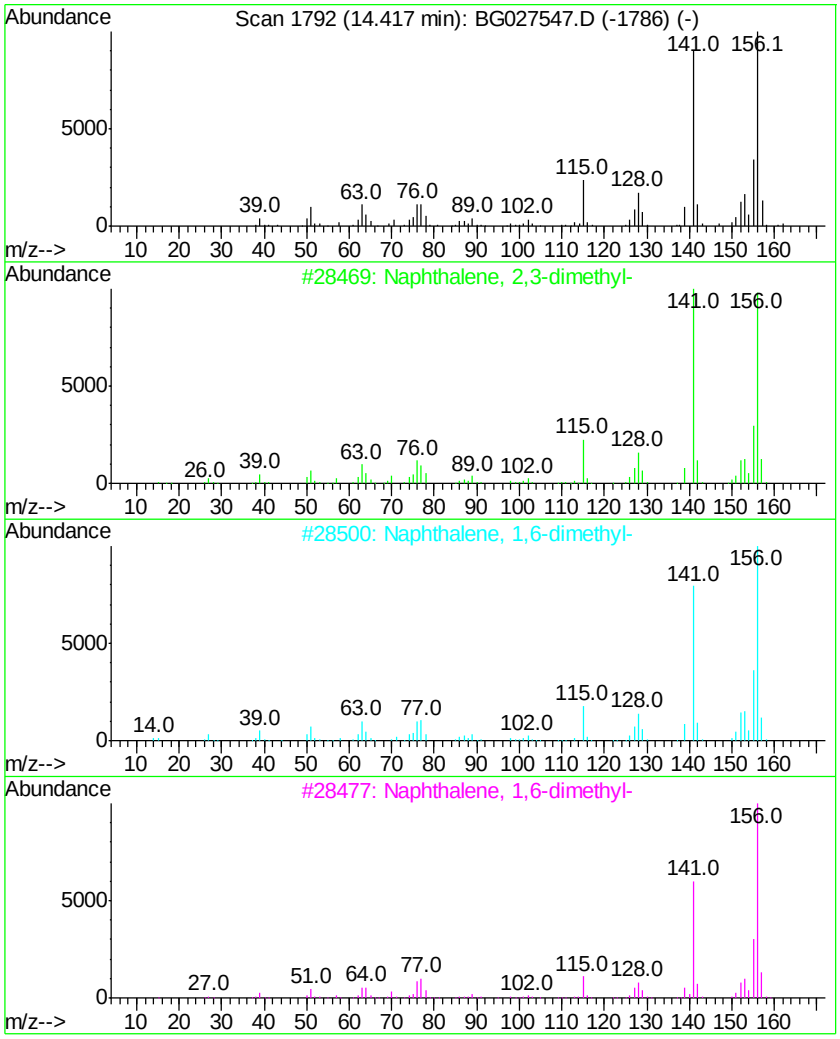
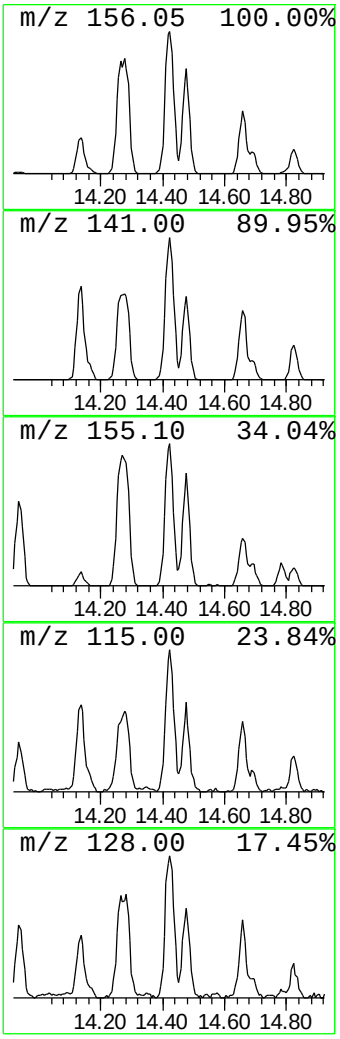
Peak Number 8 Naphthalene, 2,3-dimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.42	19.60 ng/ul	536436	Acenaphthene-d10	15.10

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, 1,6-dimethyl-	156	C12H12	000575-43-9	96
4	Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	96
5	Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	96





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Instrument : BNA_G
ClientSampleId : F5E24

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Quant Title : SVOA CALIBRATION

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

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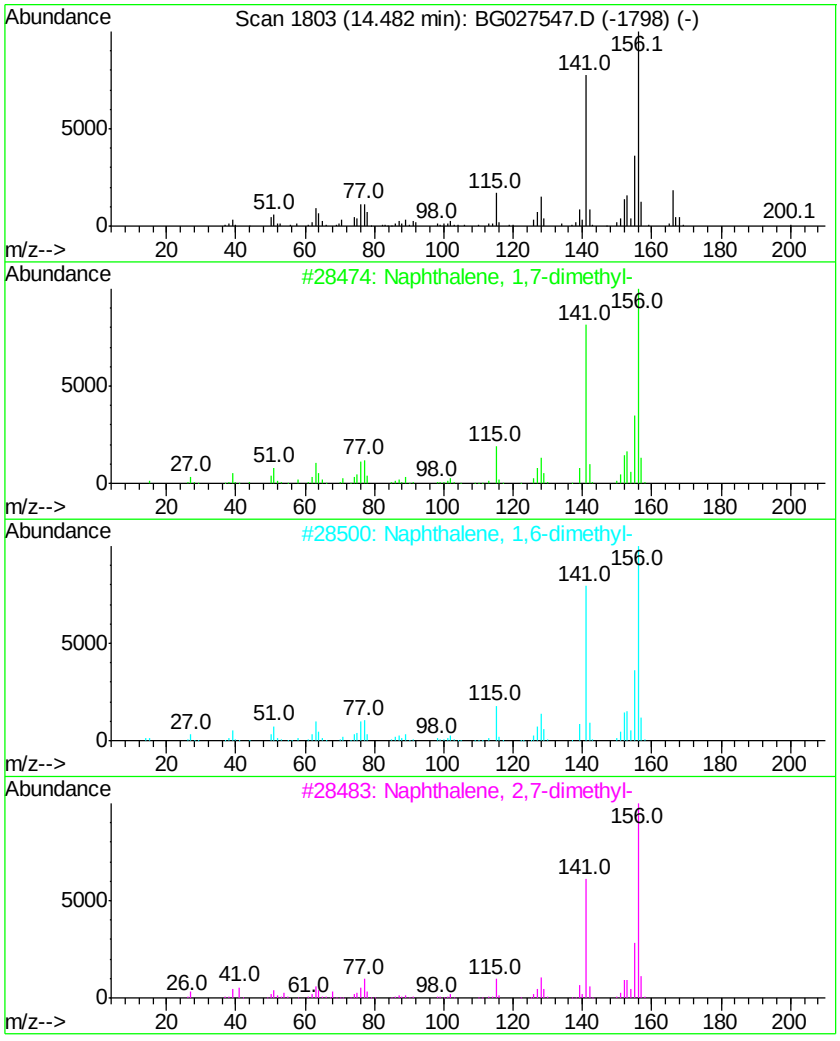
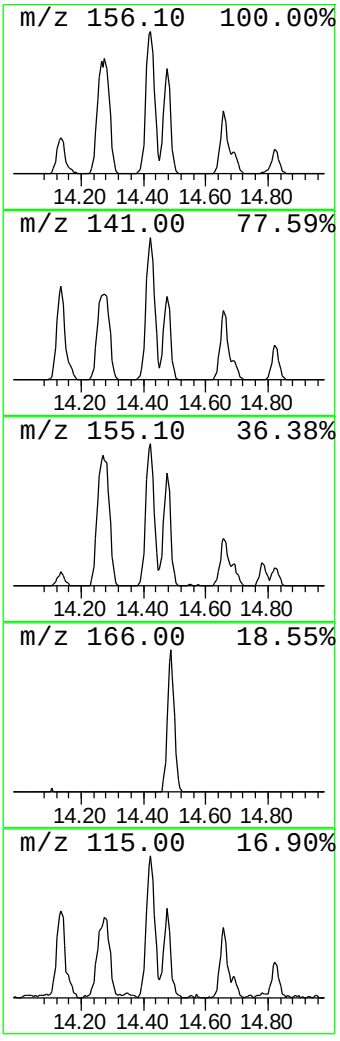
Peak Number 9 Naphthalene, 1,7-dimethyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.48	12.14 ng/ul	332211	Acenaphthene-d10	15.10

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	98
2	Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	97
3	Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	96
4	Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	96
5	Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	96





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ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : F5E24

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Quant Title : SVOA CALIBRATION

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

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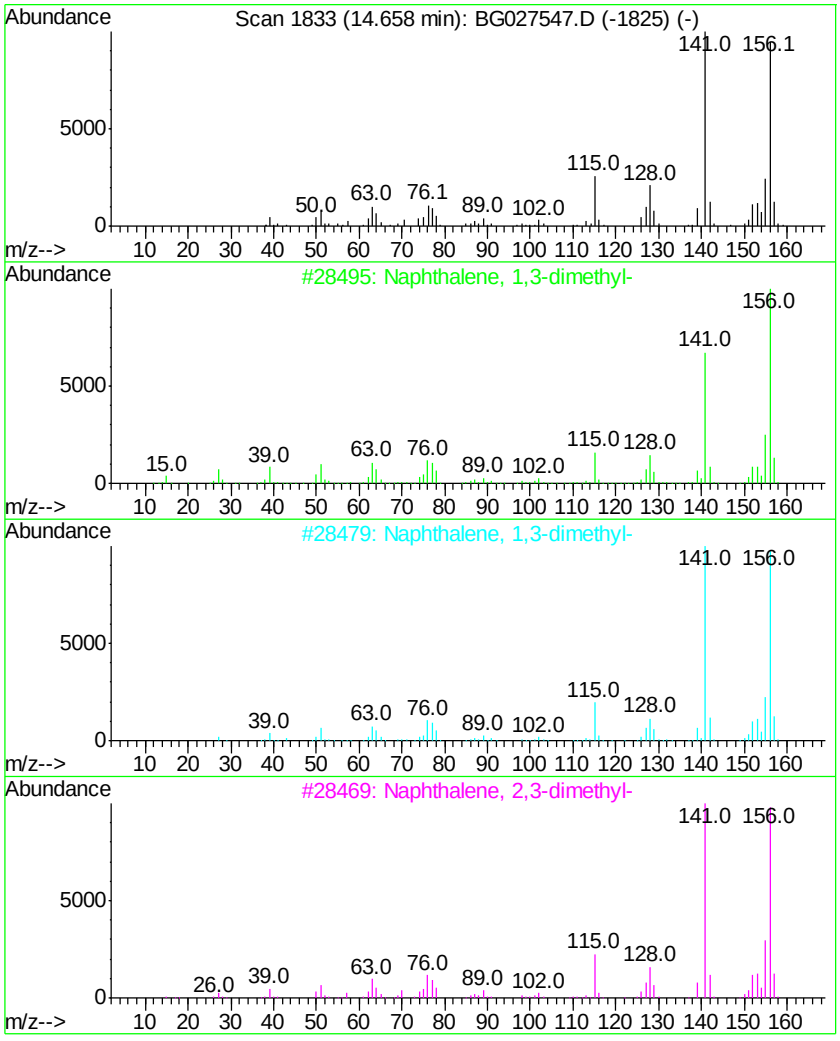
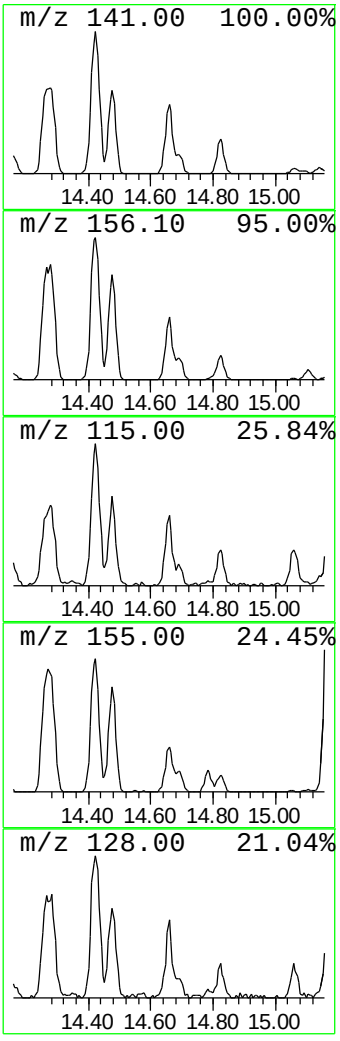
Peak Number 10 Naphthalene, 1,3-dimethyl- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.66	9.72 ng/ul	265975	Acenaphthene-d10	15.10

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	97
2	Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	97
3	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
4	Naphthalene, 1,2-dimethyl-	156	C12H12	000573-98-8	96
5	Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	96





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Sample : I3627-13 25X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : F5E24

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG061617.M
Quant Title : SVOA CALIBRATION

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

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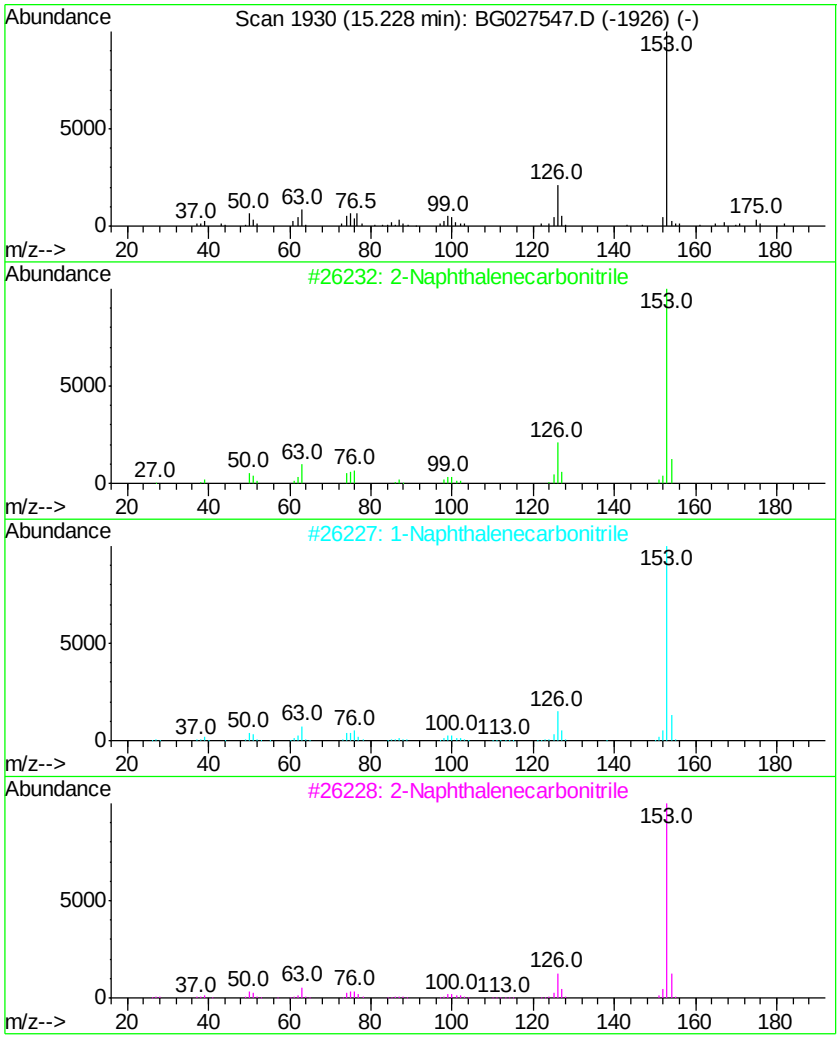
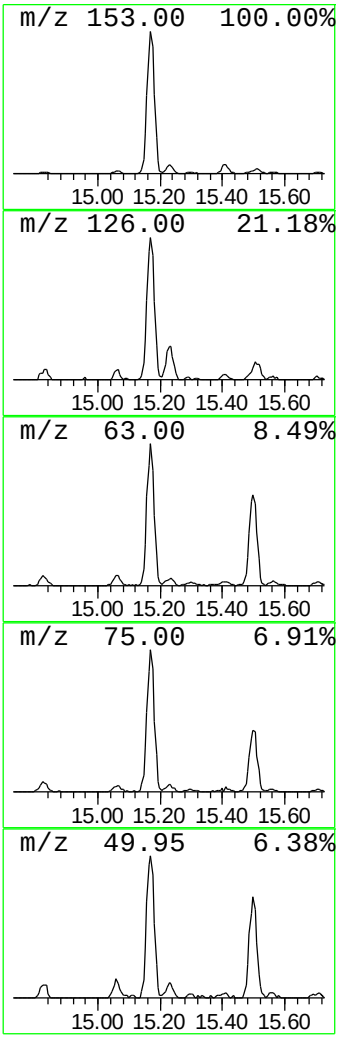
Peak Number 11 2-Naphthalenecarbonitrile Concentration Rank 45

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.23	2.94 ng/ul	80423	Acenaphthene-d10	15.10

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Naphthalenecarbonitrile	153	C11H7N	000613-46-7	91
2	1-Naphthalenecarbonitrile	153	C11H7N	000086-53-3	91
3	2-Naphthalenecarbonitrile	153	C11H7N	000613-46-7	91
4	1-Naphthalenecarbonitrile	153	C11H7N	000086-53-3	90
5	1-Naphthalenecarbonitrile	153	C11H7N	000086-53-3	87





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Sample : I3627-13 25X
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ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_G
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Quant Title : SVOA CALIBRATION

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

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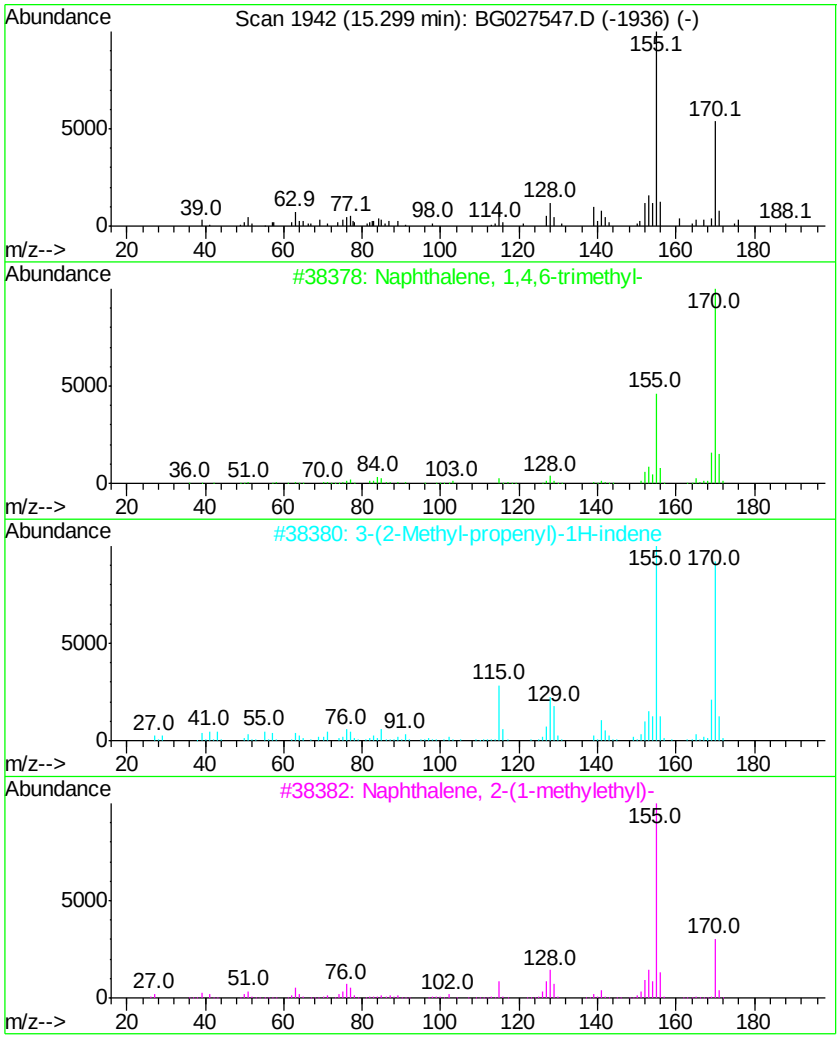
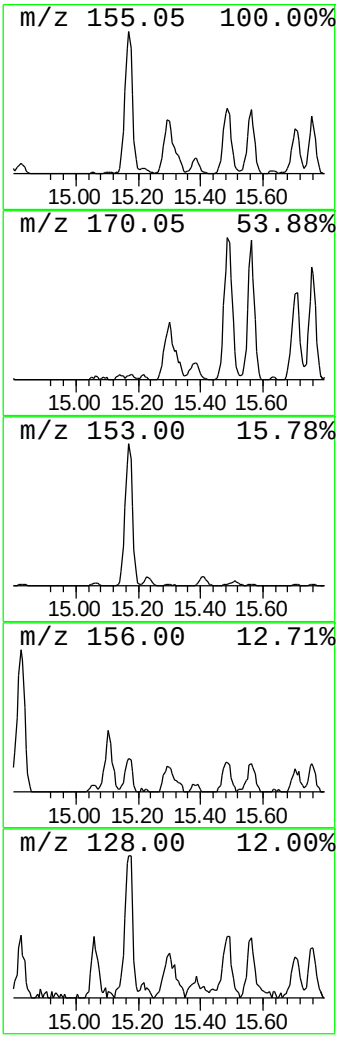
Peak Number 12 Naphthalene, 1,4,6-trimethyl- Concentration Rank 35

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.30	4.34 ng/ul	118716	Acenaphthene-d10	15.10

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,4,6-trimethyl-	170	C13H14	002131-42-2	93
2	3-(2-Methyl-propenyl)-1H-indene	170	C13H14	1000187-78-5	93
3	Naphthalene, 2-(1-methylethyl)-	170	C13H14	002027-17-0	90
4	Naphthalene, 2-(1-methylethyl)-	170	C13H14	002027-17-0	90
5	Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	87





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Sample : I3627-13 25X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : F5E24

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG061617.M
Quant Title : SVOA CALIBRATION

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

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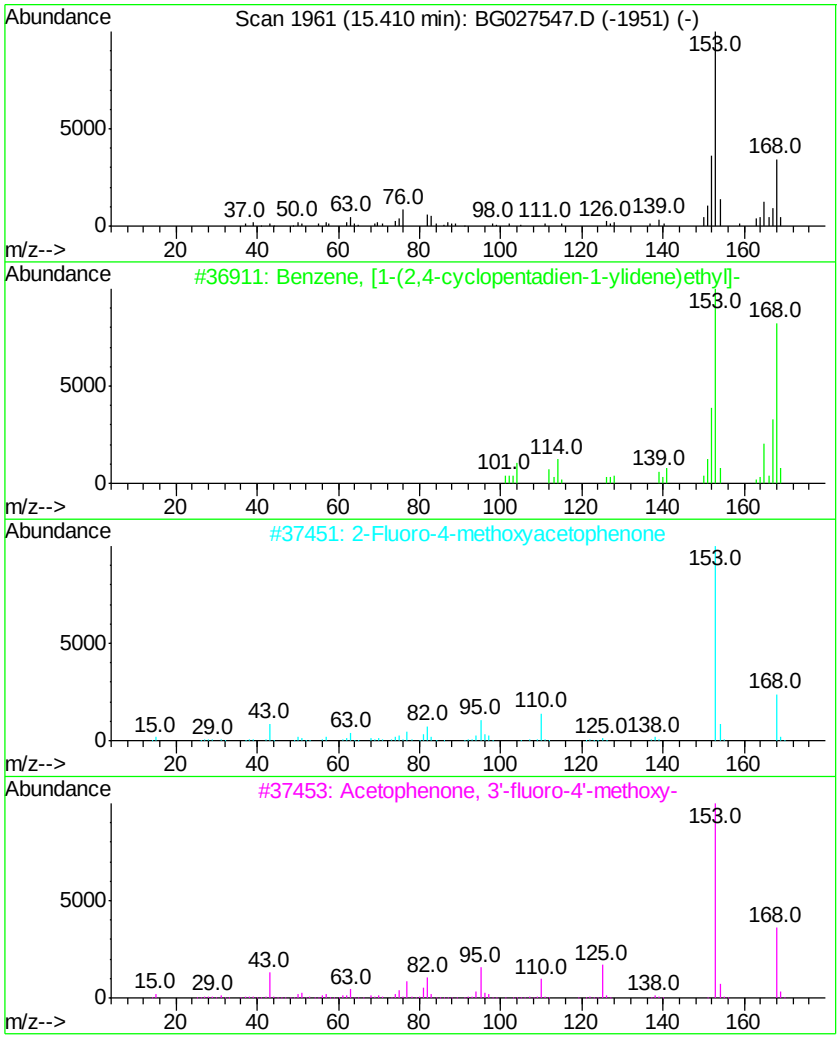
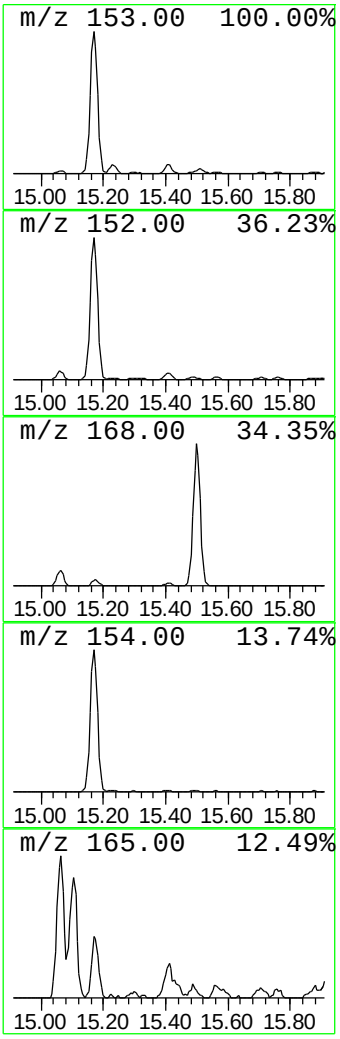
Peak Number 13 Benzene, [1-(2,4-cyclopenta... Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.41	5.63 ng/ul	154097	Acenaphthene-d10	15.10

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, [1-(2,4-cyclopentadien-...	168	C13H12	002320-32-3	56
2	2-Fluoro-4-methoxyacetophenone	168	C9H9FO2	074457-86-6	52
3	Acetophenone, 3'-fluoro-4'-methoxy-	168	C9H9FO2	000455-91-4	50
4	Ethanone, 1-(2,3,4-trihydroxyphe...	168	C8H8O4	000528-21-2	50
5	Ethanone, 1-(2,4,6-trihydroxyphe...	168	C8H8O4	000480-66-0	50





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Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : F5E24

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Quant Title : SVOA CALIBRATION

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

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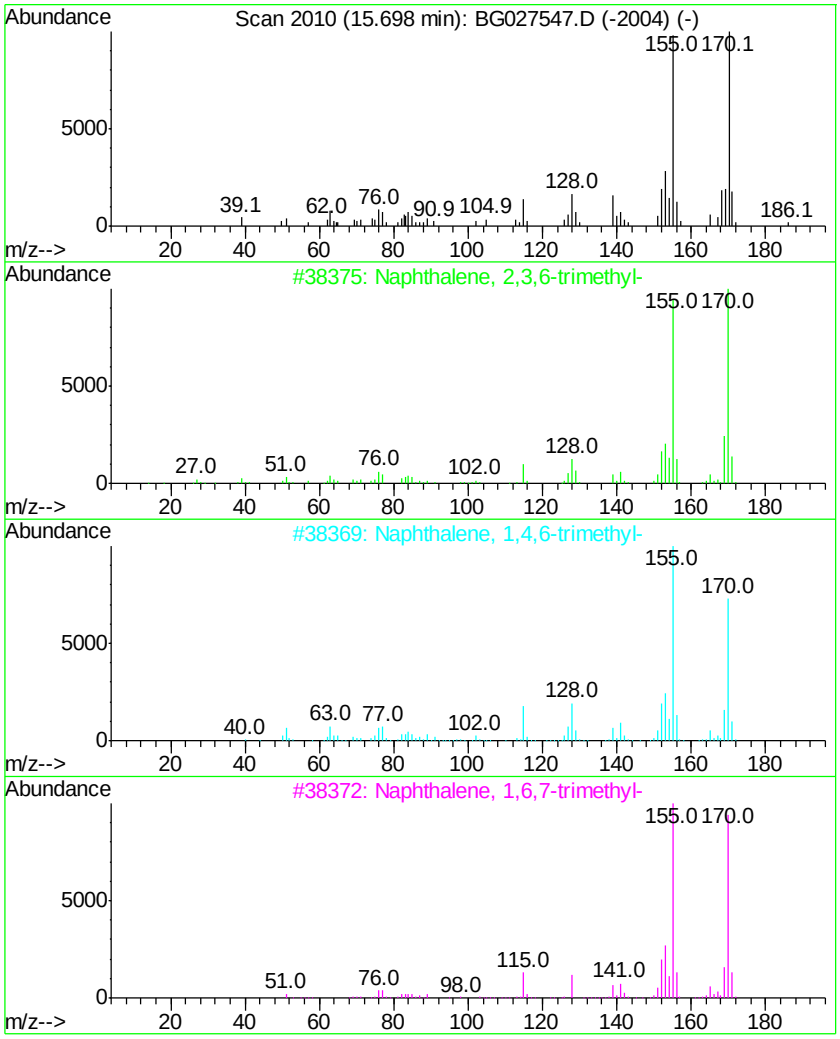
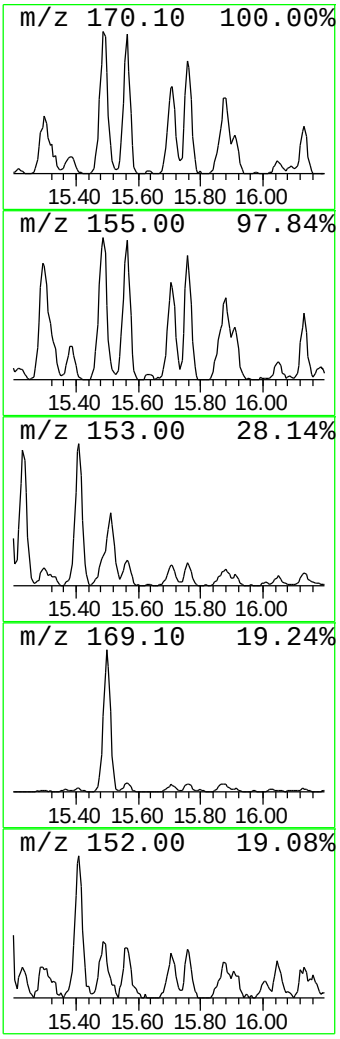
Peak Number 15 Naphthalene, 2,3,6-trimethyl- Concentration Rank 40

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.70	4.02 ng/ul	109908	Acenaphthene-d10	15.10

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	98
2	Naphthalene, 1,4,6-trimethyl-	170	C13H14	002131-42-2	97
3	Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	96
4	Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	96
5	Naphthalene, 1,4,5-trimethyl-	170	C13H14	002131-41-1	94





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Quant Title : SVOA CALIBRATION

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

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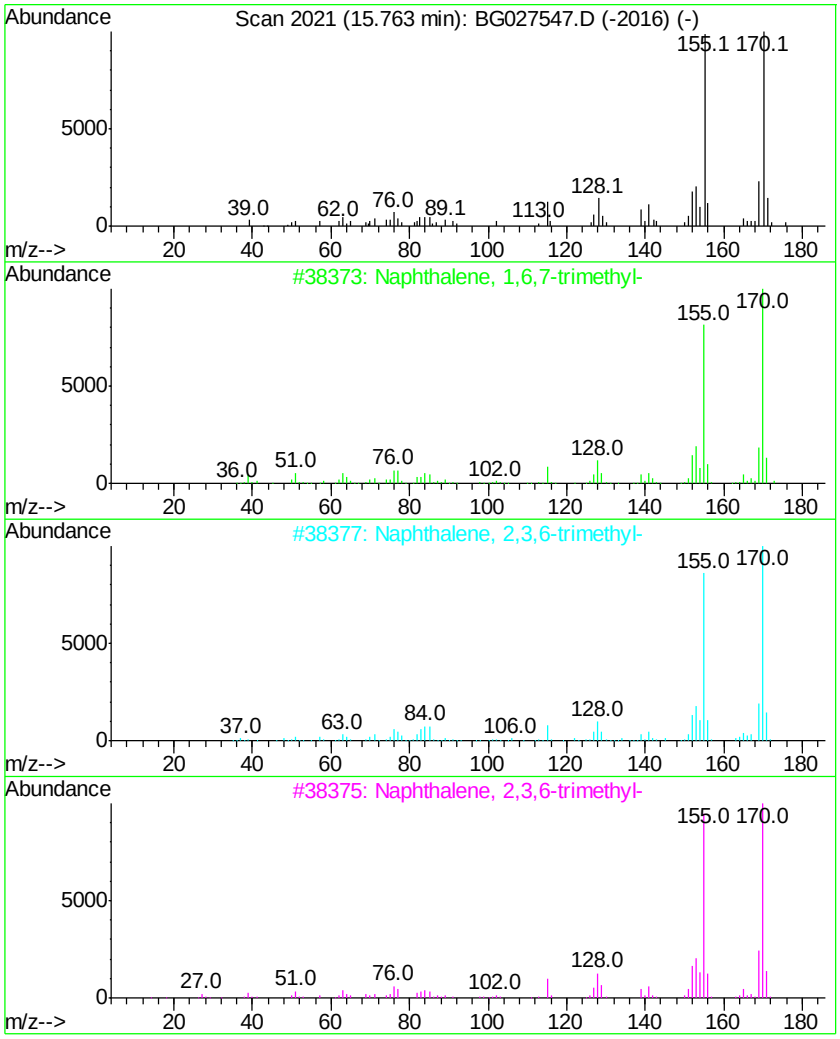
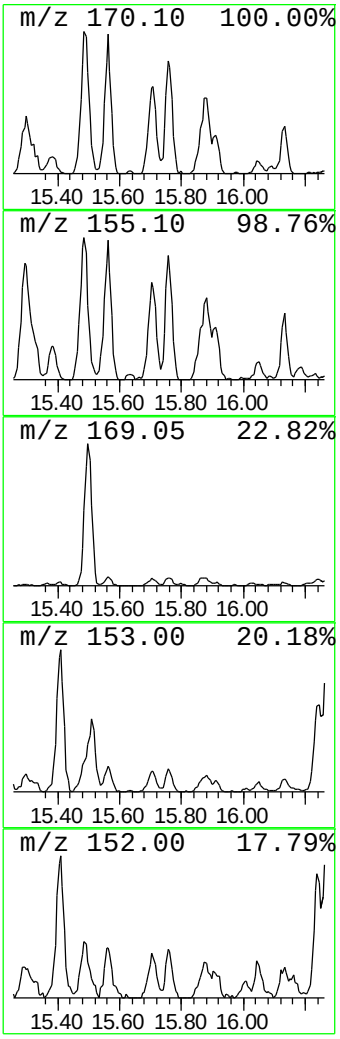
Peak Number 16 Naphthalene, 1,6,7-trimethyl- Concentration Rank 37

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.76	4.20 ng/ul	114987	Acenaphthene-d10	15.10

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	98
2	Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	97
3	Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	96
4	Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	96
5	Naphthalene, 1,4,6-trimethyl-	170	C13H14	002131-42-2	96





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Quant Title : SVOA CALIBRATION

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

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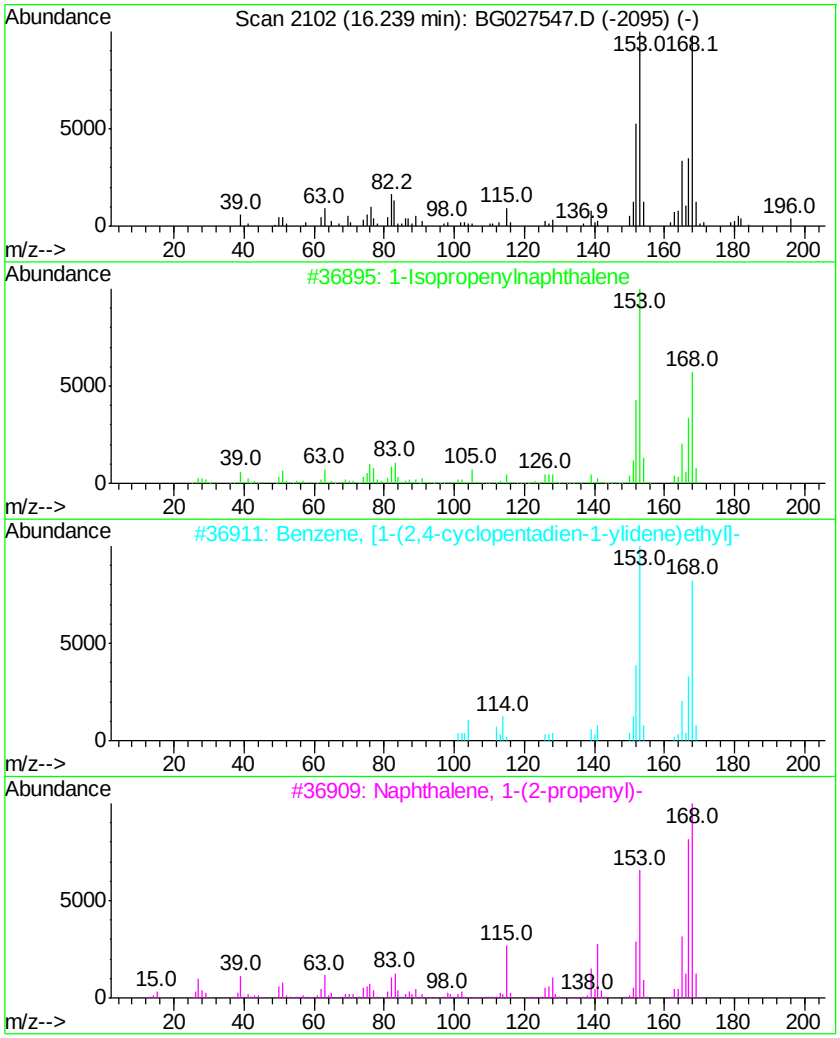
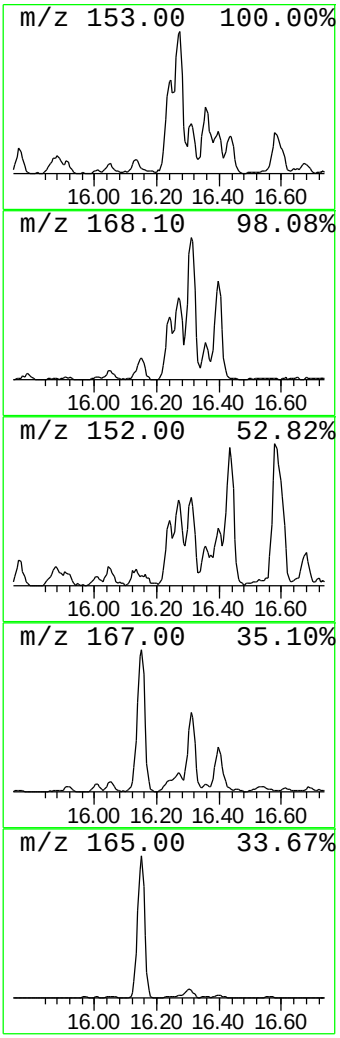
Peak Number 18 1-Isopropenylnaphthalene Concentration Rank 38

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.24	4.15 ng/ul	113587	Acenaphthene-d10	15.10

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Isopropenylnaphthalene	168	C13H12	001855-47-6	91
2	Benzene, [1-(2,4-cyclopentadien-...	168	C13H12	002320-32-3	90
3	Naphthalene, 1-(2-propenyl)-	168	C13H12	002489-86-3	83
4	1,1'-Biphenyl, 2-methyl-	168	C13H12	000643-58-3	58
5	1,1'-Biphenyl, 3-methyl-	168	C13H12	000643-93-6	43





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Quant Title : SVOA CALIBRATION

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

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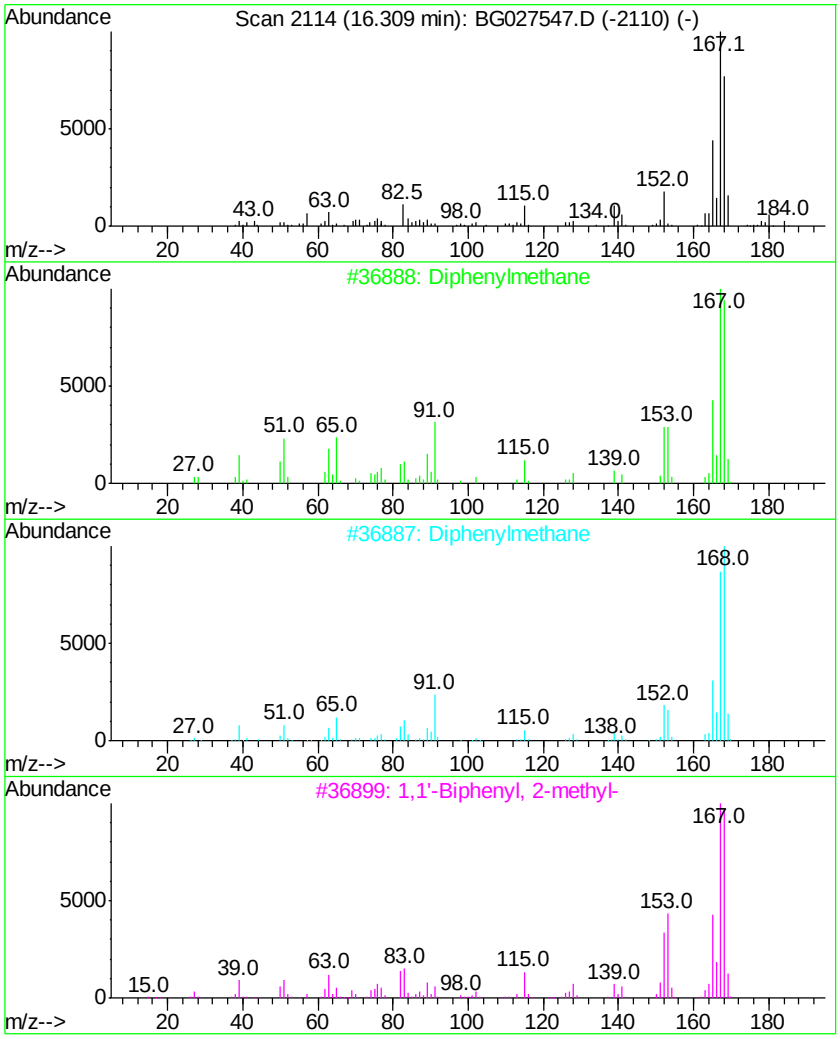
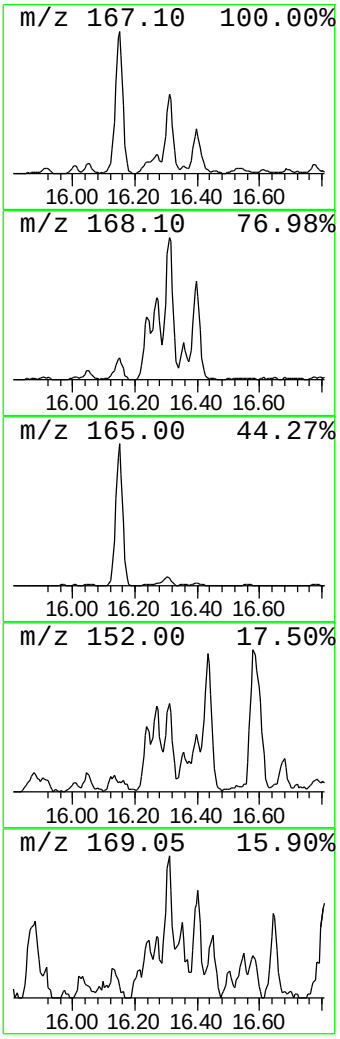
Peak Number 20 Diphenylmethane Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.31	11.90 ng/ul	325824	Acenaphthene-d10	15.10

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Diphenylmethane	168	C13H12	000101-81-5	64
2	Diphenylmethane	168	C13H12	000101-81-5	59
3	1,1'-Biphenyl, 2-methyl-	168	C13H12	000643-58-3	59
4	Fluorene, 2,4a-dihydro-	168	C13H12	059247-36-8	58
5	Naphthalene, 1-(2-propenyl)-	168	C13H12	002489-86-3	50





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Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : F5E24

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Quant Title : SVOA CALIBRATION

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

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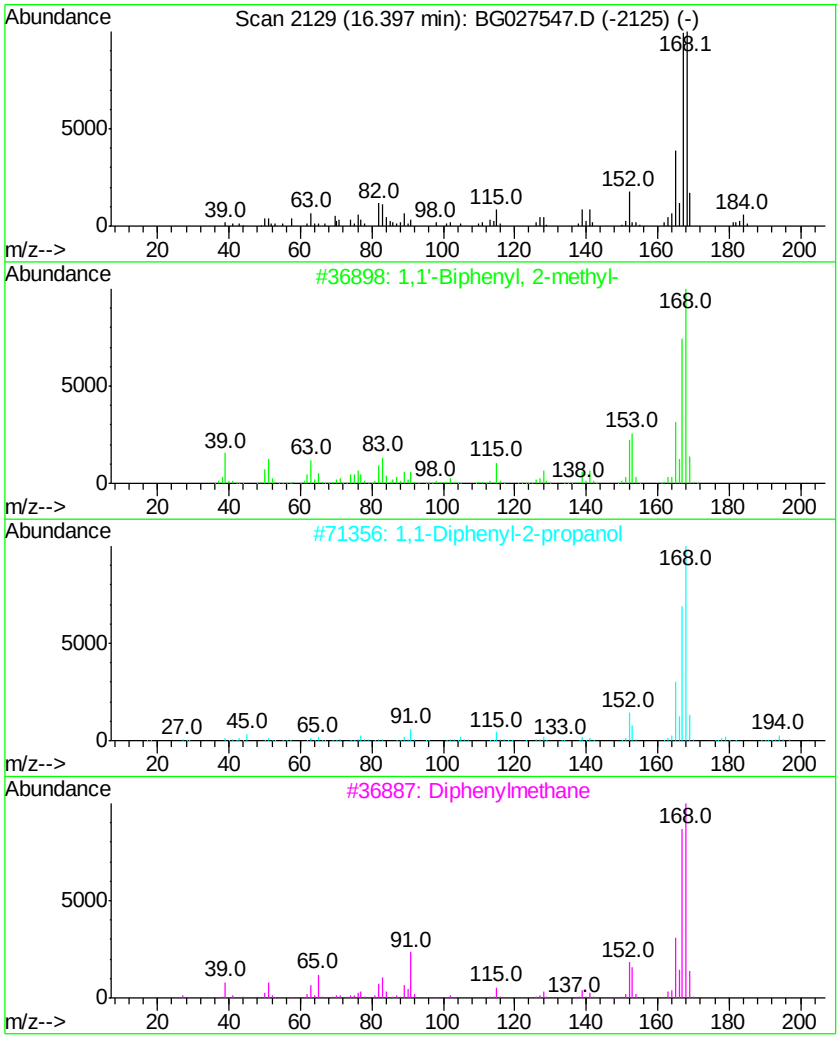
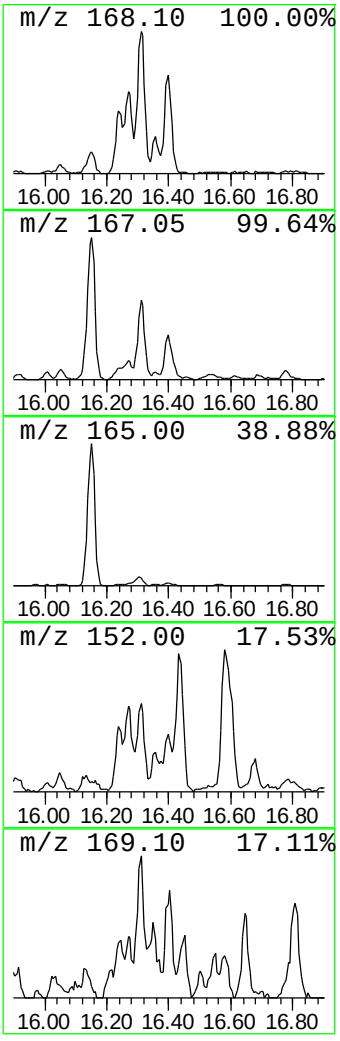
Peak Number 22 1,1'-Biphenyl, 2-methyl- Concentration Rank 28

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.40	5.62 ng/ul	153795	Acenaphthene-d10	15.10

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2-methyl-	168	C13H12	000643-58-3	90
2	1,1-Diphenyl-2-propanol	212	C15H16O	029338-49-6	83
3	Diphenylmethane	168	C13H12	000101-81-5	81
4	Diphenylmethane	168	C13H12	000101-81-5	81
5	Diphenylmethane	168	C13H12	000101-81-5	76





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Data File : BG027547.D
Acq On : 20 Jun 2017 2:56
Operator : SJ/MA
Sample : I3627-13 25X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : F5E24

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG061617.M
Quant Title : SVOA CALIBRATION

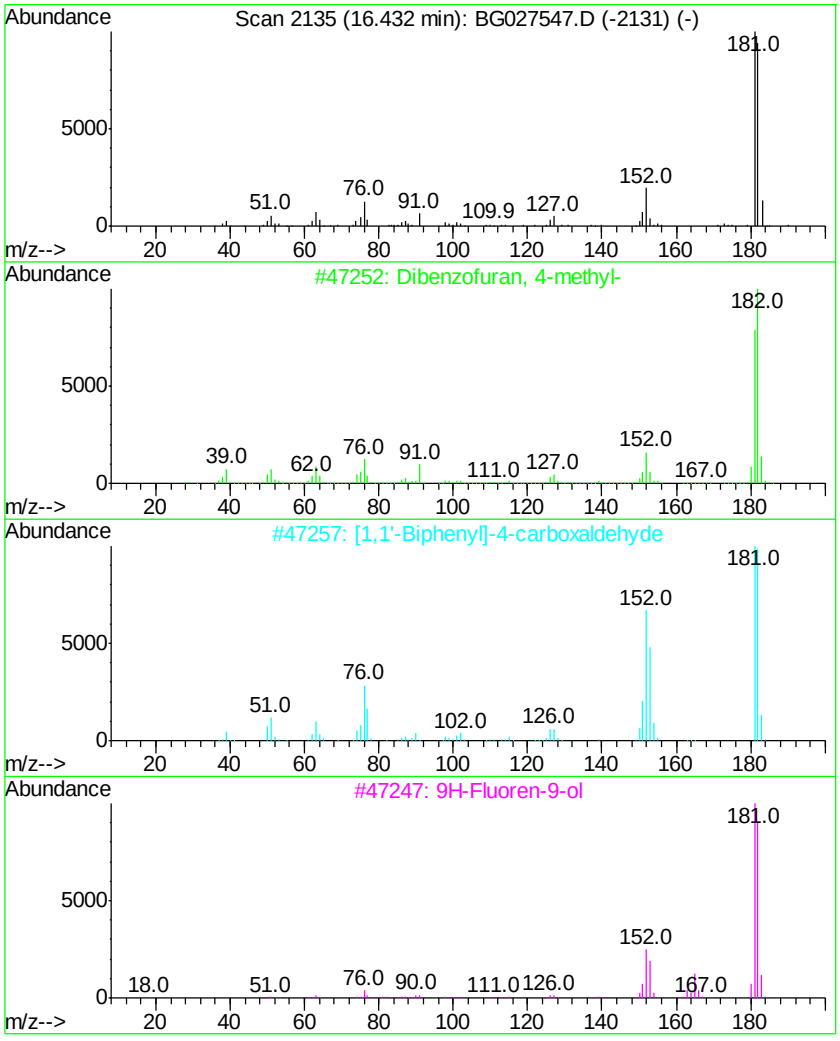
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

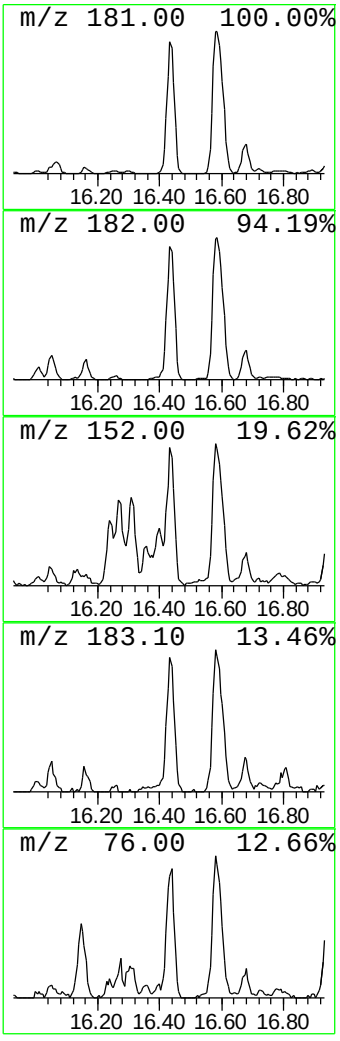
****

Peak Number 23 Dibenzofuran, 4-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.43	13.79 ng/ul	377311	Acenaphthene-d10	15.10

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dibenzofuran, 4-methyl-	182	C13H10O	007320-53-8	90
2	[1,1'-Biphenyl]-4-carboxaldehyde	182	C13H10O	003218-36-8	87
3	9H-Fluoren-9-ol	182	C13H10O	001689-64-1	86
4	[1,1'-Biphenyl]-4-carboxaldehyde	182	C13H10O	003218-36-8	83
5	[1,1'-Biphenyl]-4-carboxaldehyde	182	C13H10O	003218-36-8	78





Library Search Compound Report

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Data File : BG027547.D
Acq On : 20 Jun 2017 2:56
Operator : SJ/MA
Sample : I3627-13 25X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : F5E24

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG061617.M
Quant Title : SVOA CALIBRATION

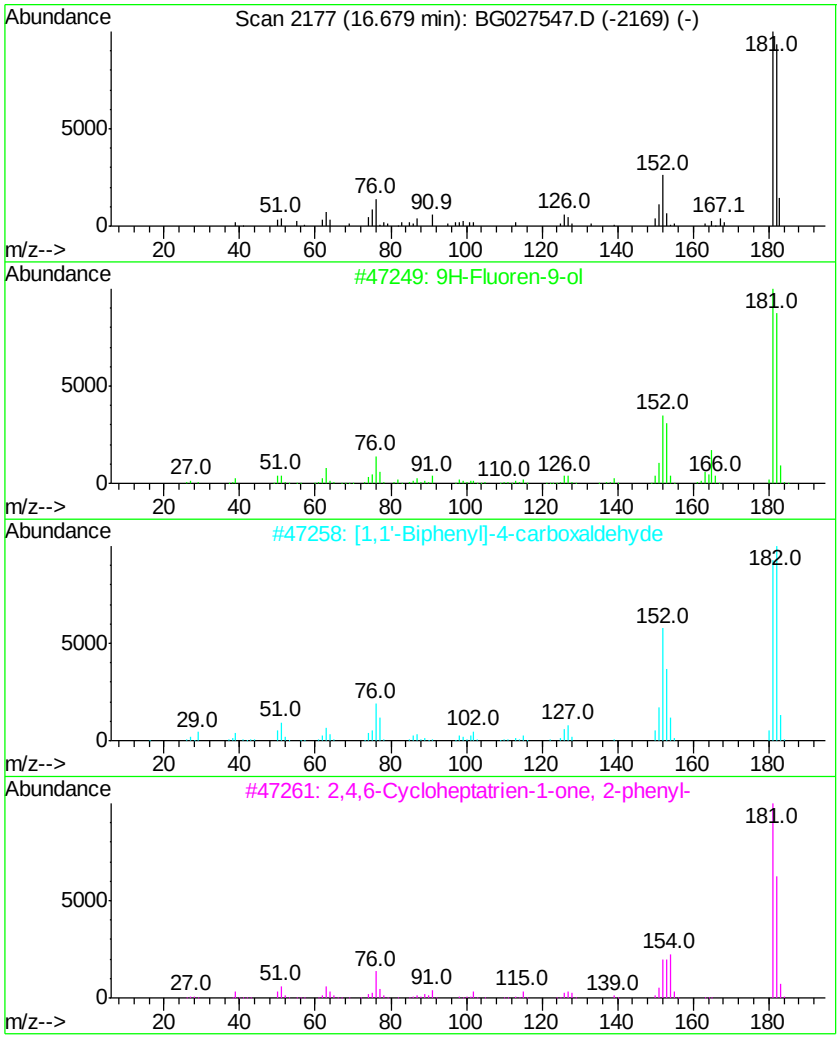
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

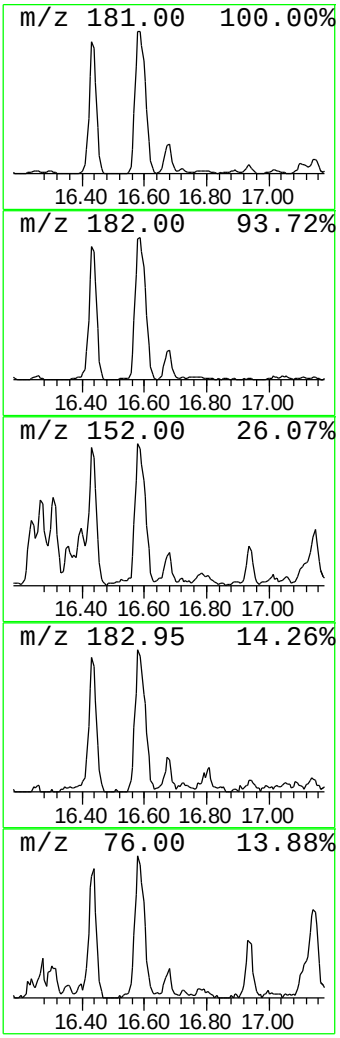
****

Peak Number 25 9H-Fluoren-9-ol Concentration Rank 48

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.68	2.81 ng/ul	100320	Phenanthrene-d10	17.85

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9H-Fluoren-9-ol	182	C13H10O	001689-64-1	91
2	[1,1'-Biphenyl]-4-carboxaldehyde	182	C13H10O	003218-36-8	91
3	2,4,6-Cycloheptatrien-1-one, 2-p...	182	C13H10O	014562-09-5	90
4	Dibenzofuran, 4-methyl-	182	C13H10O	007320-53-8	87
5	[1,1'-Biphenyl]-4-carboxaldehyde	182	C13H10O	003218-36-8	87





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Acq On : 20 Jun 2017 2:56
Operator : SJ/MA
Sample : I3627-13 25X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : F5E24

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG061617.M
Quant Title : SVOA CALIBRATION

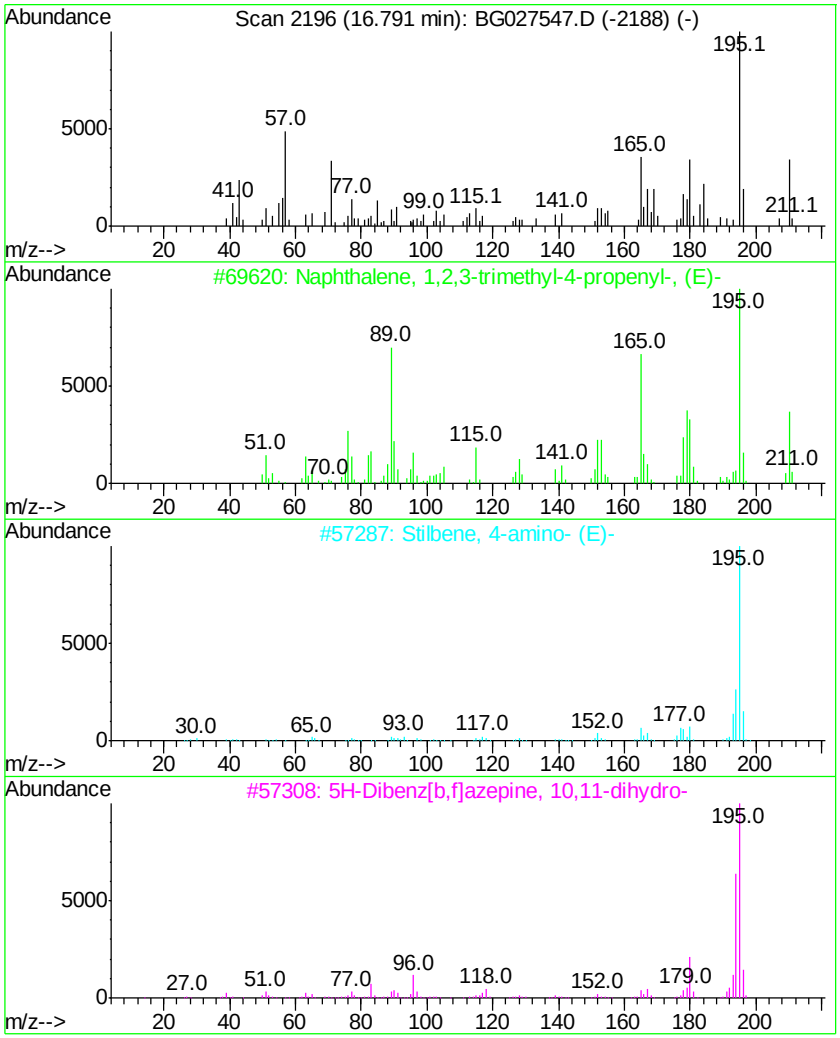
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

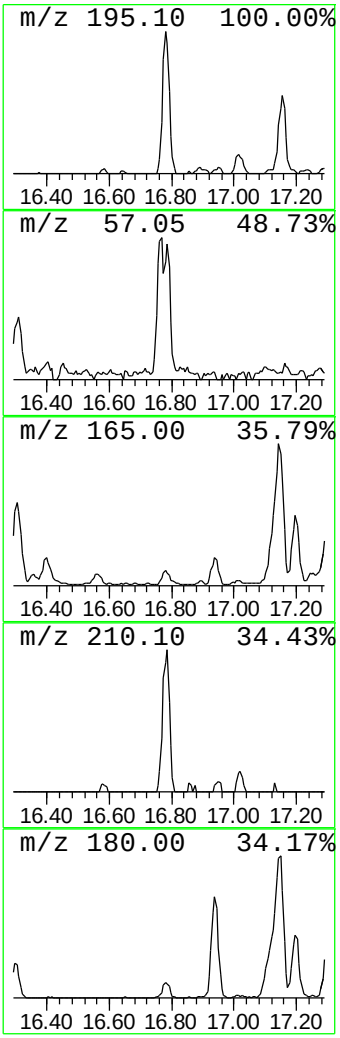
****

Peak Number 26 Naphthalene, 1,2,3-trimethy... Concentration Rank 31

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.79	5.26 ng/ul	187918	Phenanthrene-d10	17.85

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3-trimethyl-4-p...	210	C16H18	026137-53-1	64
2	Stilbene, 4-amino- (E)-	195	C14H13N	004309-66-4	58
3	5H-Dibenz[b,f]azepine, 10,11-dih...	195	C14H13N	000494-19-9	52
4	Benzenamine, 3-(2-phenylethenyl)...	195	C14H13N	014064-82-5	49
5	(4-Acetylphenyl)phenylmethane	210	C15H14O	000782-92-3	49





Library Search Compound Report

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Data File : BG027547.D
Acq On : 20 Jun 2017 2:56
Operator : SJ/MA
Sample : I3627-13 25X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : F5E24

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG061617.M
Quant Title : SVOA CALIBRATION

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

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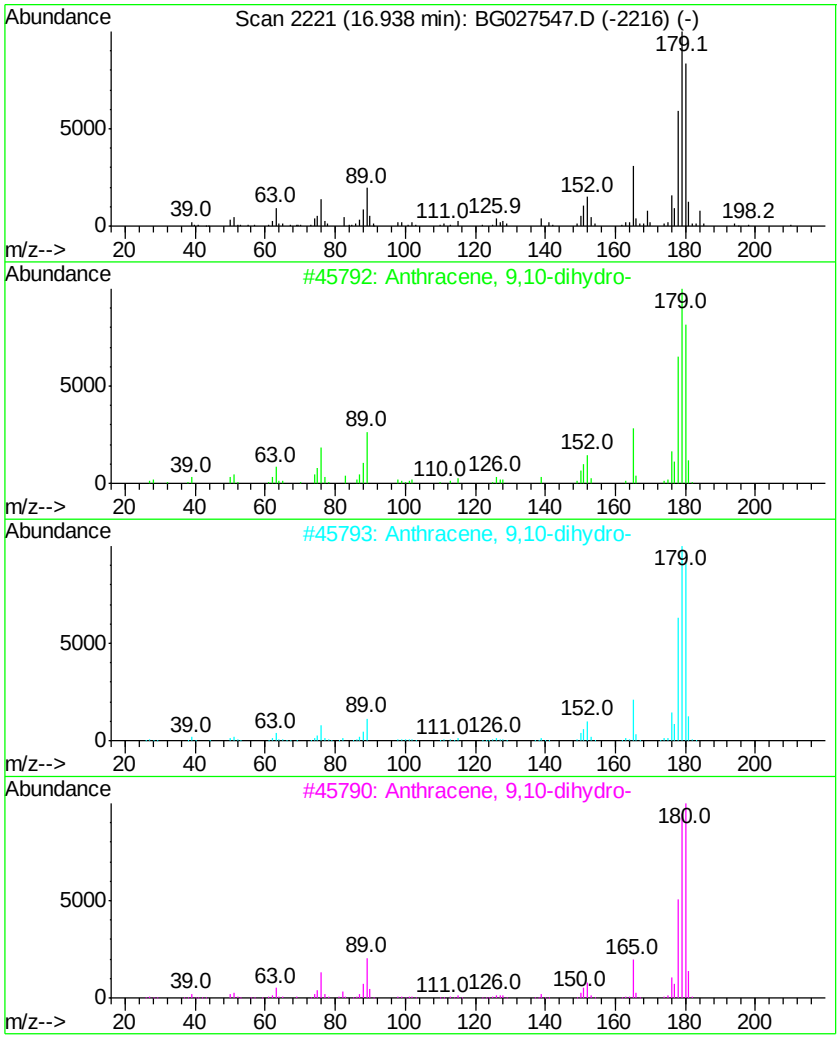
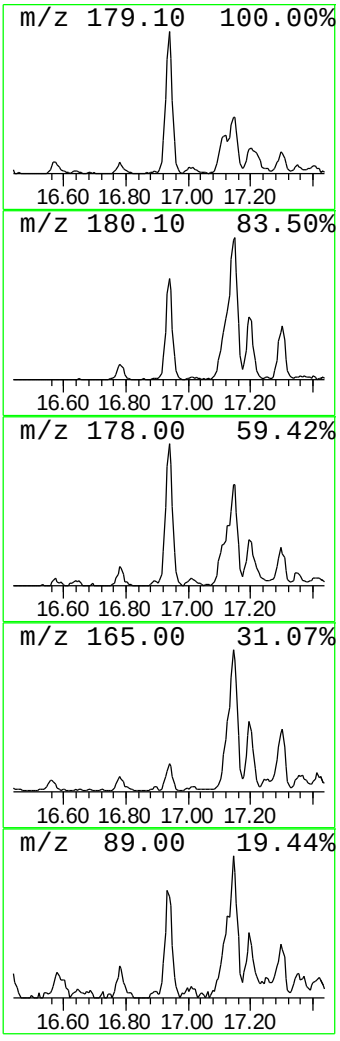
Peak Number 27 Anthracene, 9,10-dihydro- Concentration Rank 30

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.94	5.39 ng/ul	192680	Phenanthrene-d10	17.85

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Anthracene, 9,10-dihydro-	180	C14H12	000613-31-0	98
2	Anthracene, 9,10-dihydro-	180	C14H12	000613-31-0	97
3	Anthracene, 9,10-dihydro-	180	C14H12	000613-31-0	97
4	Phenanthrene, 9,10-dihydro-	180	C14H12	000776-35-2	96
5	Phenanthrene, 9,10-dihydro-	180	C14H12	000776-35-2	96





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Acq On : 20 Jun 2017 2:56
Operator : SJ/MA
Sample : I3627-13 25X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : F5E24

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG061617.M
Quant Title : SVOA CALIBRATION

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

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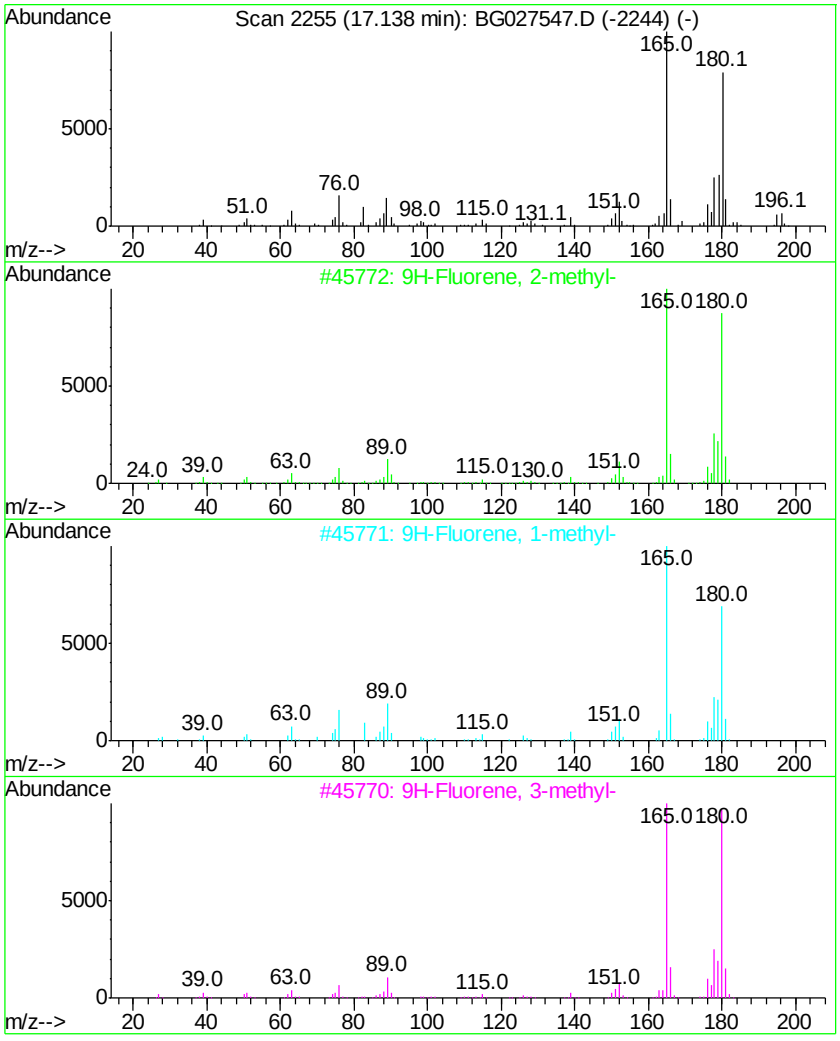
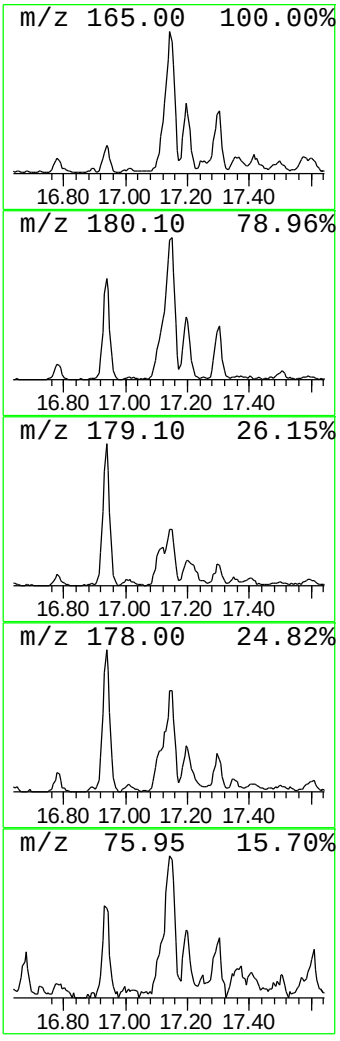
Peak Number 28 9H-Fluorene, 2-methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.14	12.87 ng/ul	459913	Phenanthrene-d10	17.85

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9H-Fluorene, 2-methyl-	180	C14H12	001430-97-3	96
2	9H-Fluorene, 1-methyl-	180	C14H12	001730-37-6	96
3	9H-Fluorene, 3-methyl-	180	C14H12	002523-39-9	93
4	9H-Fluorene, 2-methyl-	180	C14H12	001430-97-3	92
5	9H-Fluorene, 1-methyl-	180	C14H12	001730-37-6	92





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Data File : BG027547.D
Acq On : 20 Jun 2017 2:56
Operator : SJ/MA
Sample : I3627-13 25X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : F5E24

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG061617.M
Quant Title : SVOA CALIBRATION

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

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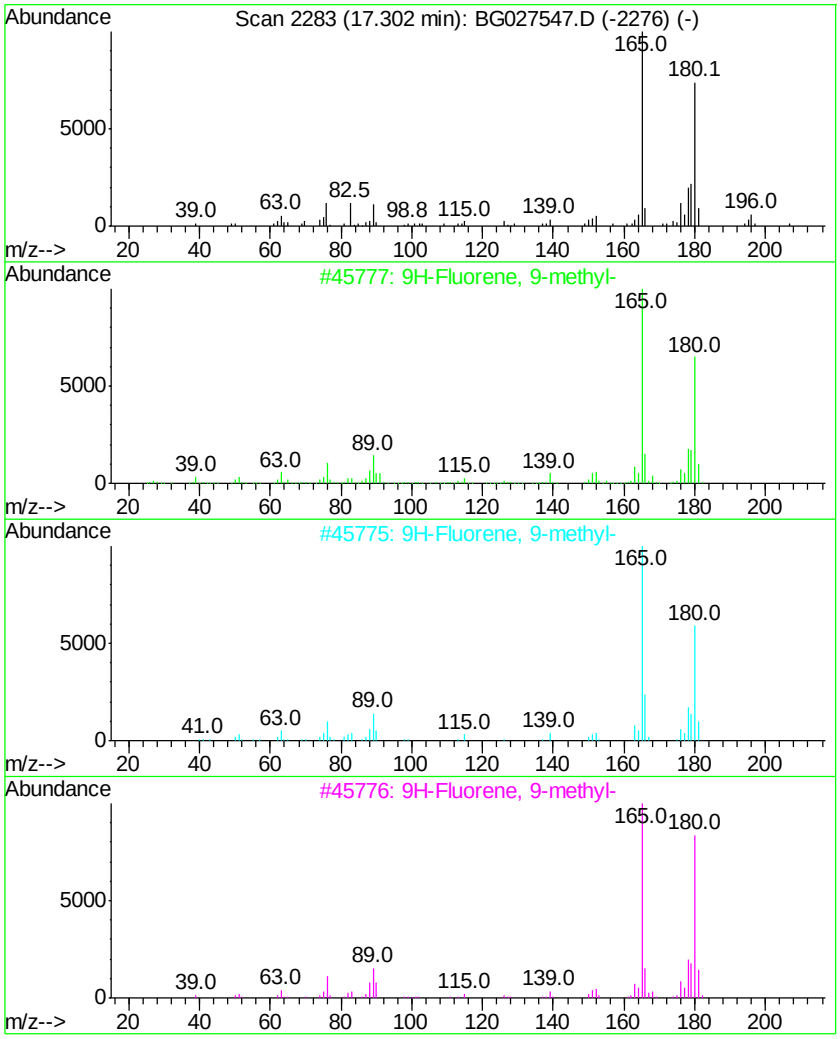
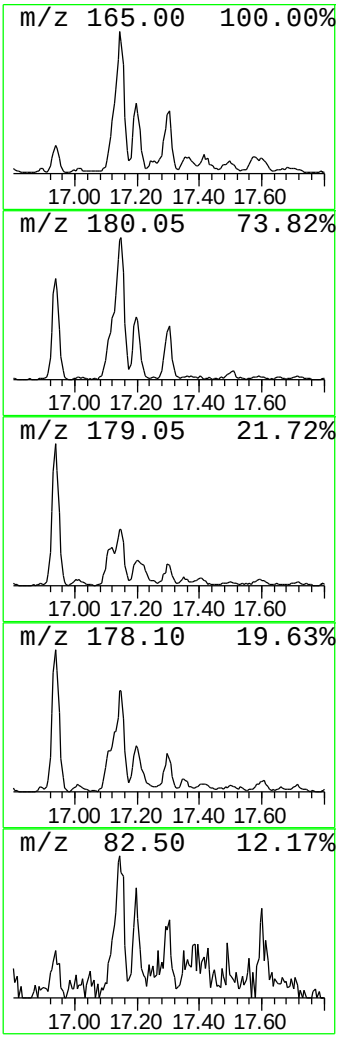
Peak Number 30 9H-Fluorene, 9-methyl- Concentration Rank 39

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.30	4.02 ng/ul	143643	Phenanthrene-d10	17.85

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9H-Fluorene, 9-methyl-	180	C14H12	002523-37-7	72
2	9H-Fluorene, 9-methyl-	180	C14H12	002523-37-7	68
3	9H-Fluorene, 9-methyl-	180	C14H12	002523-37-7	68
4	9H-Fluorene, 1-methyl-	180	C14H12	001730-37-6	64
5	9H-Fluorene, 1-methyl-	180	C14H12	001730-37-6	64





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Acq On : 20 Jun 2017 2:56
Operator : SJ/MA
Sample : I3627-13 25X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : F5E24

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG061617.M
Quant Title : SVOA CALIBRATION

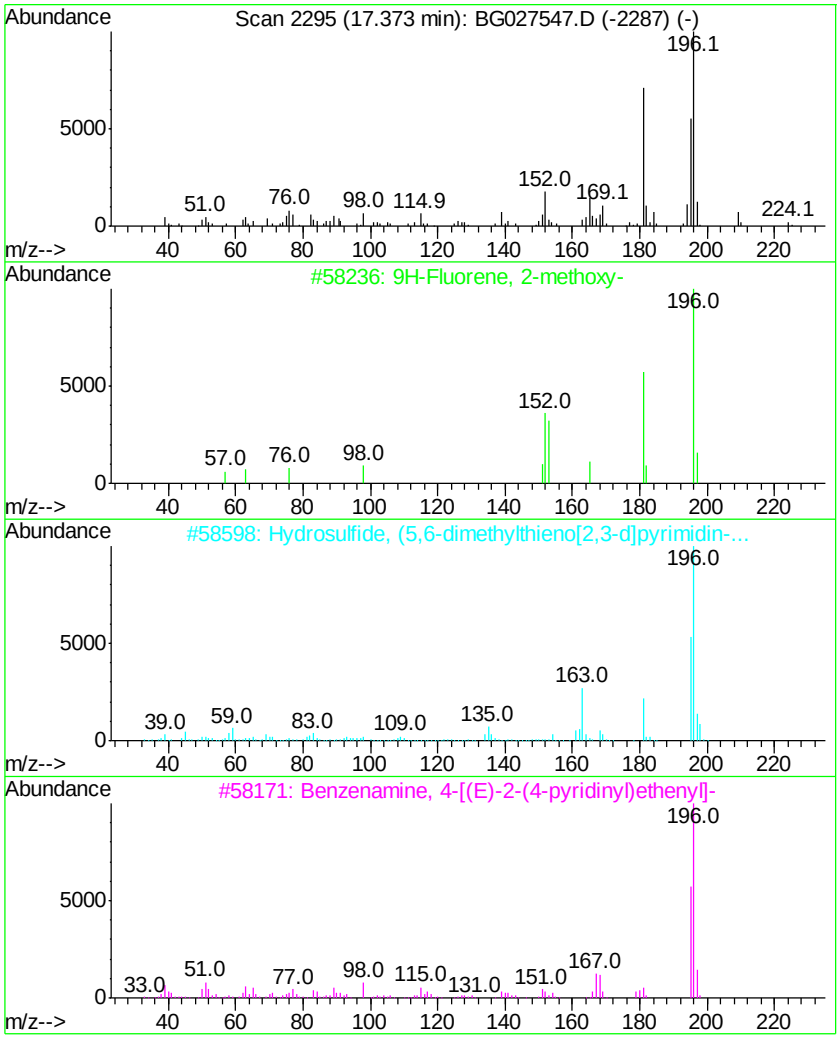
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

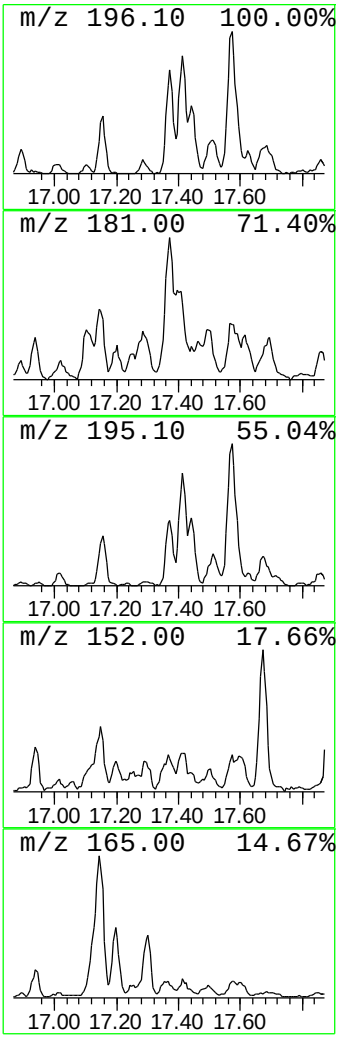
****

Peak Number 31 9H-Fluorene, 2-methoxy- Concentration Rank 32

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.37	5.13 ng/ul	183358	Phenanthrene-d10	17.85

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9H-Fluorene, 2-methoxy-	196	C14H12O	002523-46-8	50
2	Hydrosulfide, (5,6-dimethylthien...	196	C8H8N2S2	1000362-41-8	50
3	Benzenamine, 4-[(E)-2-(4-pyridin...	196	C13H12N2	1000351-61-4	49
4	Phenol, 3-(2-phenylethenyl)-, (E)-	196	C14H12O	017861-18-6	47
5	Phenol, 4-(2-phenylethenyl)-, (E)-	196	C14H12O	006554-98-9	38





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Operator : SJ/MA
Sample : I3627-13 25X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : F5E24

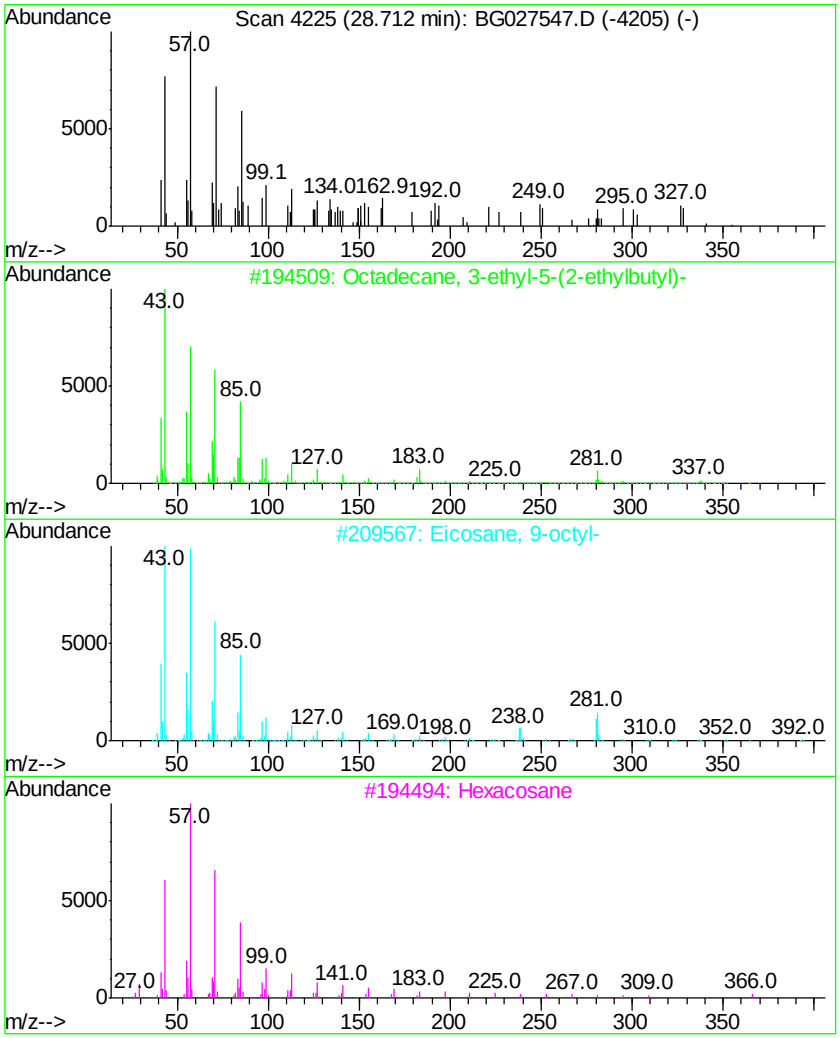
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG061617.M
Quant Title : SVOA CALIBRATION

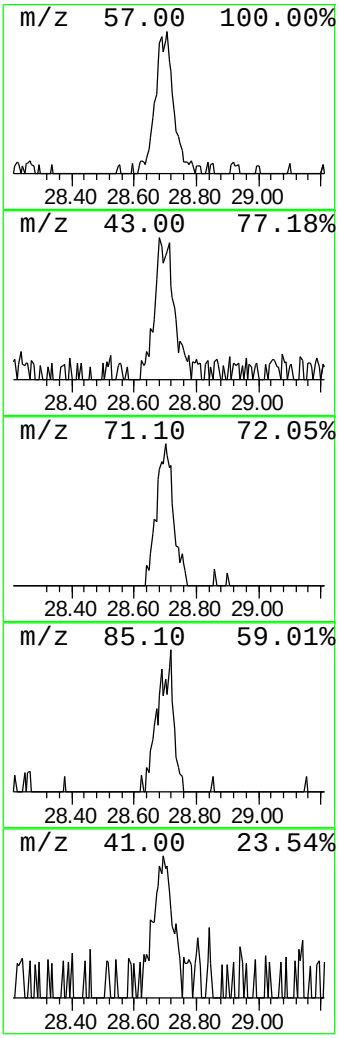
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 55 (DEL) Alkane: Straight-Chai... Concentration Rank 55

R.T.	EstConc	Area	Relative to ISTD	R.T.
28.71	2.04 ng/ul	76776	Perylene-d12	25.88

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecane, 3-ethyl-5-(2-ethylbu...	366	C26H54	055282-12-7	58
2	Eicosane, 9-octyl-	394	C28H58	013475-77-9	50
3	Hexacosane	366	C26H54	000630-01-3	49
4	Heptacosane	380	C27H56	000593-49-7	49
5	Tetracosane, 11-decyl-	479	C34H70	055429-84-0	49





Tentatively Identified Compound (LSC) summary

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG061917\
Data File : BG027547.D
Acq On : 20 Jun 2017 2:56
Operator : SJ/MA
Sample : I3627-13 25X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : F5E24

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG061617.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
Indane	8.89	14.8	ng/ul	172871	1	8.52	233477	20.0
Indene	9.06	12.7	ng/ul	147882	1	8.52	233477	20.0
Benzene, 2-etheny...	10.72	6.1	ng/ul	121713	2	11.34	396337	20.0
Benzo[b]thiophene	11.51	11.7	ng/ul	231821	2	11.34	396337	20.0
Naphthalene, 2-et...	14.14	8.8	ng/ul	240019	3	15.10	547400	20.0
Naphthalene, 2,6-...	14.28	20.7	ng/ul	566615	3	15.10	547400	20.0
Naphthalene, 2,3-...	14.42	19.6	ng/ul	536436	3	15.10	547400	20.0
Naphthalene, 1,7-...	14.48	12.1	ng/ul	332211	3	15.10	547400	20.0
Naphthalene, 1,3-...	14.66	9.7	ng/ul	265975	3	15.10	547400	20.0
2-Naphthalenecarb...	15.23	2.9	ng/ul	80423	3	15.10	547400	20.0
Naphthalene, 1,4,...	15.30	4.3	ng/ul	118716	3	15.10	547400	20.0
Benzene, [1-(2,4-...	15.41	5.6	ng/ul	154097	3	15.10	547400	20.0
Naphthalene, 2,3,...	15.70	4.0	ng/ul	109908	3	15.10	547400	20.0
Naphthalene, 1,6,...	15.76	4.2	ng/ul	114987	3	15.10	547400	20.0
1-Isopropenylnaph...	16.24	4.2	ng/ul	113587	3	15.10	547400	20.0
Diphenylmethane	16.31	11.9	ng/ul	325824	3	15.10	547400	20.0
1,1'-Biphenyl, 2-...	16.40	5.6	ng/ul	153795	3	15.10	547400	20.0
Dibenzofuran, 4-m...	16.43	13.8	ng/ul	377311	3	15.10	547400	20.0
9H-Fluoren-9-ol	16.68	2.8	ng/ul	100320	4	17.85	714738	20.0
Naphthalene, 1,2,...	16.79	5.3	ng/ul	187918	4	17.85	714738	20.0
Anthracene, 9,10-...	16.94	5.4	ng/ul	192680	4	17.85	714738	20.0
9H-Fluorene, 2-me...	17.14	12.9	ng/ul	459913	4	17.85	714738	20.0
9H-Fluorene, 9-me...	17.30	4.0	ng/ul	143643	4	17.85	714738	20.0
9H-Fluorene, 2-me...	17.37	5.1	ng/ul	183358	4	17.85	714738	20.0
(DEL) Alkane: Str...	28.71	2.0	ng/ul	76776	6	25.88	751067	20.0