LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG080217\
Data File : BG028136.D
Acq On : 3 Aug 2017 11:33
Operator : SJ/JU
Sample : I4405-07
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : JJ2X4

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-BG080217.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	3.852	41	45	55	rVB5	5678	13553	0.78%	0.052%
2	4.710	186	191	197	rBV5	8156	13298	0.76%	0.051%
3	5.115	250	260	263	rVB4	5868	12242	0.70%	0.047%
4	5.879	383	390	397	rVB	21669	34765	2.00%	0.134%
5	6.355	461	471	483	rBV	108489	190583	10.95%	0.737%
6	6.849	545	555	561	rVB3	9488	13965	0.80%	0.054%
7	7.507	659	667	683	rVB4	189897	430054	24.71%	1.664%
8	7.677	689	696	706	rBV	84387	150241	8.63%	0.581%
9	7.783	708	714	720	rVB4	10305	17071	0.98%	0.066%
10	7.900	726	734	742	rBV	111195	203329	11.68%	0.787%
11	8.365	806	813	827	rVB	138269	260997	14.99%	1.010%
12	8.605	847	854	862	rVB4	5200	11447	0.66%	0.044%
13	9.058	923	931	943	rVV2	162167	307113	17.64%	1.188%
14	9.187	945	953	964	rBV	101506	185944	10.68%	0.719%
15	9.534	1004	1012	1021	rVV2	128795	252228	14.49%	0.976%
16	9.810	1052	1059	1068	rVB5	6102	12031	0.69%	0.047%
17	10.256	1126	1135	1143	rBV	116527	214043	12.30%	0.828%
18	10.803	1209	1228	1241	rBV	171092	352490	20.25%	1.363%
19	11.185	1286	1293	1307	rVB	240899	458295	26.33%	1.773%
20	11.320	1307	1316	1325	rBV4	23646	52517	3.02%	0.203%
21	11.567	1352	1358	1369	rVB2	15144	29265	1.68%	0.113%
22	12.101	1439	1449	1456	rBV5	7604	23585	1.35%	0.091%
23	12.389	1492	1498	1505	rBV5	5119	10866	0.62%	0.042%
24	12.930	1583	1590	1606	rBV2	73109	219815	12.63%	0.850%
25	13.394	1663	1669	1677	rBV3	11177	26706	1.53%	0.103%
26	13.459	1677	1680	1688	rVB7	5643	10830	0.62%	0.042%
27	13.564	1693	1698	1705	rVB6	5873	12597	0.72%	0.049%
28	13.782	1724	1735	1745	rBV6	25102	58767	3.38%	0.227%
29	14.093	1780	1788	1800	rBV2	31666	69746	4.01%	0.270%
30	14.363	1825	1834	1838	rBV	388600	614326	35.29%	2.376%
31	14.405	1838	1841	1852	rVB	105573	159820	9.18%	0.618%
32	14.669	1877	1886	1895	rVV	406608	674113	38.73%	2.608%
33	14.828	1909	1913	1921	rBV7	17580	28065	1.61%	0.109%
34	14.898	1921	1925	1928	rVV4	8086	10615	0.61%	0.041%

LSC Area Percent Report

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Data File : BG028136.D
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Operator : SJ/JU
Sample : I4405-07
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : JJ2X4

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-BG080217.M
Title : SVOA CALIBRATION

35	14.974	1928	1938	1946	rVB2	407894	707160	40.62%	2.735%
36	15.115	1955	1962	1966	rBV4	20847	36059	2.07%	0.139%
37	15.168	1966	1971	1980	rVV3	71032	156974	9.02%	0.607%
38	15.286	1986	1991	1993	rBV5	13304	19902	1.14%	0.077%
39	15.356	1999	2003	2014	rVB10	14496	44289	2.54%	0.171%
40	15.533	2028	2033	2040	rBV3	35762	55988	3.22%	0.217%
41	15.773	2070	2074	2080	rVB3	18257	26804	1.54%	0.104%
42	15.961	2099	2106	2112	rVB	567773	884411	50.81%	3.421%
43	16.020	2112	2116	2119	rBV5	10399	16281	0.94%	0.063%
44	16.067	2120	2124	2135	rVB2	151677	276667	15.89%	1.070%
45	16.179	2136	2143	2145	rVV6	11455	23568	1.35%	0.091%
46	16.226	2145	2151	2156	rVB7	22030	46536	2.67%	0.180%
47	16.314	2162	2166	2170	rBV7	9825	17337	1.00%	0.067%
48	16.367	2170	2175	2183	rVV5	40829	73650	4.23%	0.285%
49	16.437	2183	2187	2192	rVV2	16513	23436	1.35%	0.091%
50	16.502	2192	2198	2208	rVB	354555	507762	29.17%	1.964%
51	16.590	2208	2213	2216	rBV6	8730	12522	0.72%	0.048%
52	16.637	2216	2221	2228	rBV3	64101	107202	6.16%	0.415%
53	16.725	2228	2236	2246	rVV	49224	104018	5.98%	0.402%
54	16.819	2246	2252	2258	rVB10	9803	23265	1.34%	0.090%
55	16.972	2273	2278	2280	rBV4	29202	46749	2.69%	0.181%
56	17.048	2281	2291	2307	rVB4	229303	615290	35.35%	2.380%
57	17.178	2309	2313	2323	rVB9	31134	66907	3.84%	0.259%
58	17.254	2323	2326	2330	rBV2	14724	24110	1.39%	0.093%
59	17.307	2331	2335	2338	rVB	29685	34346	1.97%	0.133%
60	17.348	2339	2342	2345	rBV	13683	12206	0.70%	0.047%
61	17.430	2352	2356	2359	rVB3	17720	20997	1.21%	0.081%
62	17.477	2360	2364	2367	rVV5	29308	50402	2.90%	0.195%
63	17.518	2367	2371	2378	rVB2	79687	114504	6.58%	0.443%
64	17.624	2385	2389	2391	rBV2	13803	17552	1.01%	0.068%
65	17.665	2391	2396	2400	rVV2	191367	285047	16.37%	1.103%
66	17.724	2400	2406	2410	rVV	520108	854817	49.11%	3.307%
67	17.765	2410	2413	2417	rVV	112823	160201	9.20%	0.620%
68	17.818	2417	2422	2440	rVB2	586386	1083041	62.22%	4.189%
69	18.030	2453	2458	2461	rBV7	9517	20574	1.18%	0.080%
70	18.112	2462	2472	2484	rVV2	143848	292398	16.80%	1.131%
71	18.212	2484	2489	2500	rVB2	113037	303745	17.45%	1.175%

LSC Area Percent Report

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Sample : I4405-07
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : JJ2X4

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-BG080217.M
Title : SVOA CALIBRATION

72	18.329	2504	2509	2513	rBV2	60090	86095	4.95%	0.333%
73	18.412	2516	2523	2530	rVB3	87019	218556	12.56%	0.845%
74	18.506	2535	2539	2543	rBV7	13724	20451	1.17%	0.079%
75	18.576	2545	2551	2567	rBV6	83466	251948	14.47%	0.975%
76	18.776	2581	2585	2593	rBV10	25764	48629	2.79%	0.188%
77	18.852	2595	2598	2607	rVB10	14669	22713	1.30%	0.088%
78	19.081	2631	2637	2641	rBV	112589	182675	10.49%	0.707%
79	19.122	2641	2644	2646	rBV2	45390	54417	3.13%	0.210%
80	19.510	2707	2710	2715	rVB6	20597	29541	1.70%	0.114%
81	19.575	2718	2721	2727	rBV5	24515	39023	2.24%	0.151%
82	19.698	2738	2742	2744	rBV5	20691	33358	1.92%	0.129%
83	19.857	2761	2769	2779	rVB2	177826	412107	23.67%	1.594%
84	20.098	2803	2810	2816	rVB2	770417	1155162	66.36%	4.468%
85	20.339	2848	2851	2855	rVB	44878	54209	3.11%	0.210%
86	20.585	2889	2893	2898	rVB3	53689	76561	4.40%	0.296%
87	21.050	2967	2972	2976	rBV4	55752	108671	6.24%	0.420%
88	21.179	2989	2994	3000	rBV	254907	366821	21.07%	1.419%
89	21.238	3001	3004	3011	rVB	95799	143696	8.25%	0.556%
90	21.355	3020	3024	3030	rBV5	110383	190319	10.93%	0.736%
91	21.608	3062	3067	3069	rBV	272579	408817	23.49%	1.581%
92	21.643	3069	3073	3080	rVB	889089	1330997	76.46%	5.149%
93	22.054	3136	3143	3156	rVB	512749	1006884	57.84%	3.895%
94	22.278	3175	3181	3189	rVB	1062912	1692721	97.24%	6.548%
95	22.795	3265	3269	3276	rVB2	106606	187080	10.75%	0.724%
96	23.300	3351	3355	3364	rVB2	79773	163012	9.36%	0.631%
97	25.327	3692	3700	3714	rVB2	298333	905572	52.02%	3.503%
98	25.574	3731	3742	3754	rVB2	464944	1488920	85.53%	5.759%
99	27.201	4009	4019	4032	rVB3	322003	1130266	64.93%	4.372%
100	27.595	4073	4086	4103	rVB	473307	1740746	100.00%	6.734%

Sum of corrected areas: 25852006

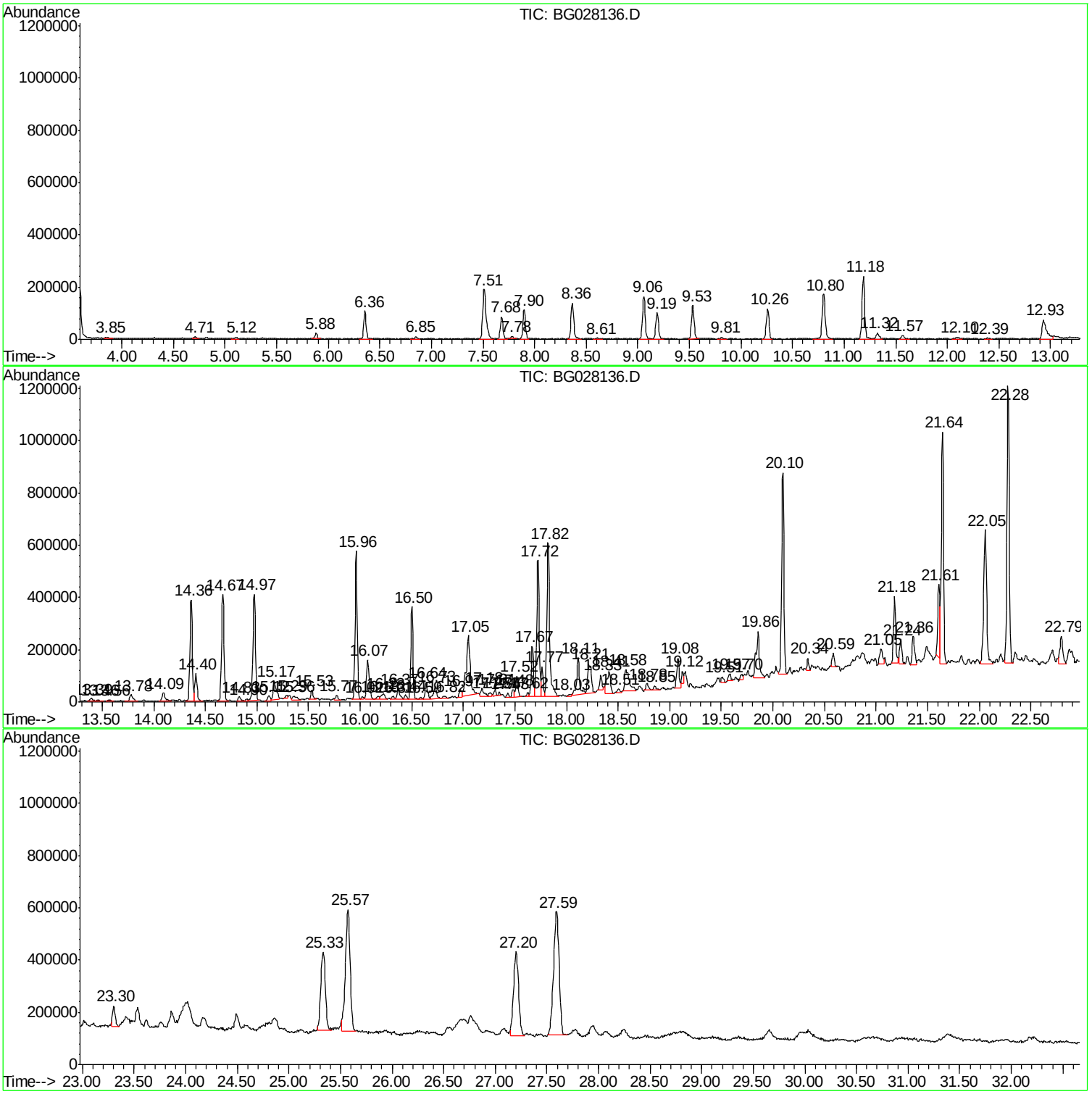
LSC Report - Integrated Chromatogram

Data Path : Z:\HPCHEM1\BNA G\DATA\BG080217\
Data File : BG028136.D
Acq On : 3 Aug 2017 11:33
Operator : SJ/JU
Sample : I4405-07
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : JJ2X4

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG080217.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\HPCHEM1\BNA G\DATA\BG080217\
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ALS Vial : 35 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : JJ2X4

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG080217.M
Quant Title : SVOA CALIBRATION

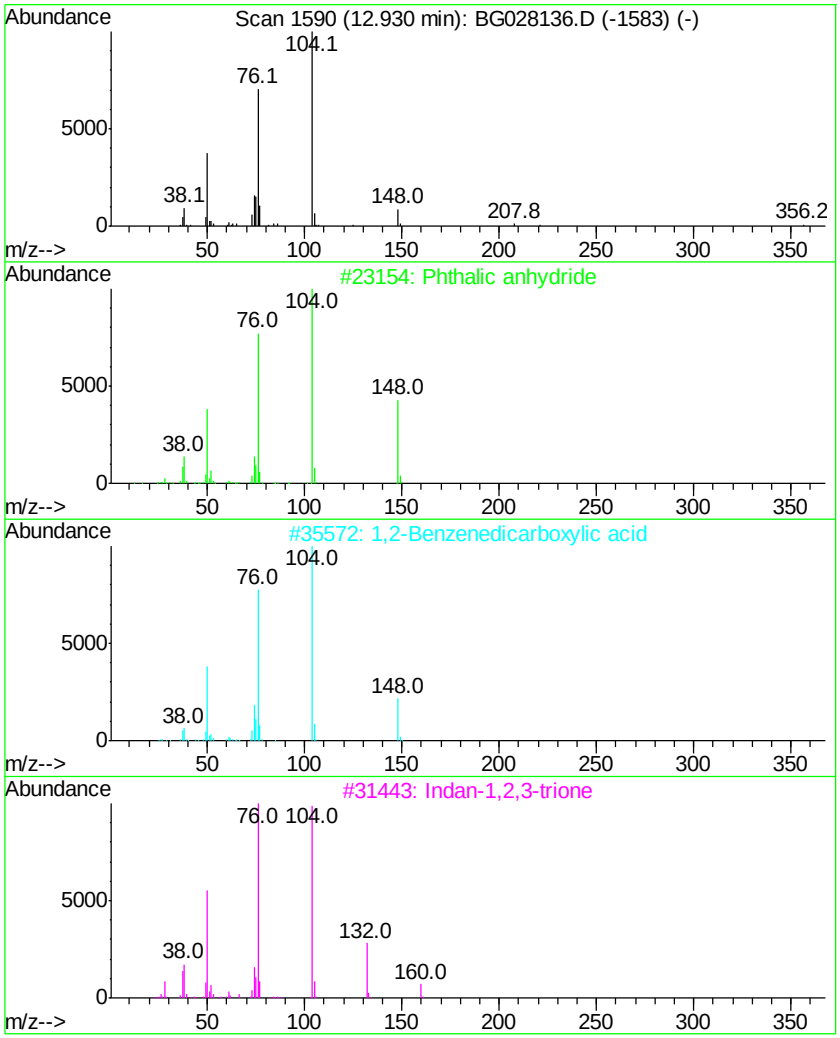
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

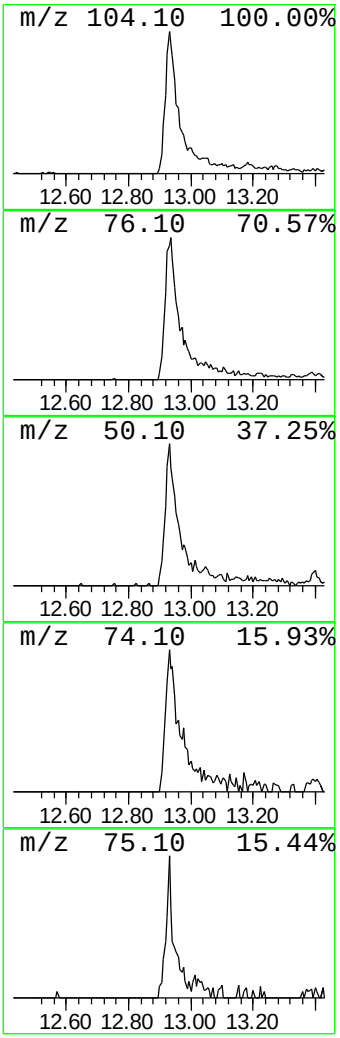
****

Peak Number 3 Phthalic anhydride Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.93	9.59 ng/ul	219815	Naphthalene-d8	11.18

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phthalic anhydride	148	C8H4O3	000085-44-9	91
2	1,2-Benzenedicarboxylic acid	166	C8H6O4	000088-99-3	90
3	Indan-1,2,3-trione	160	C9H4O3	000938-24-9	83
4	Phthalamic acid	165	C8H7NO3	000088-97-1	72
5	2-Sulfobenzoic acid cyclic anhyd...	184	C7H4O4S	000081-08-3	72





Data Path : Z:\HPCHEM1\BNA G\DATA\BG080217\
Data File : BG028136.D
Acq On : 3 Aug 2017 11:33
Operator : SJ/JU
Sample : I4405-07
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : JJ2X4

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG080217.M
Quant Title : SVOA CALIBRATION

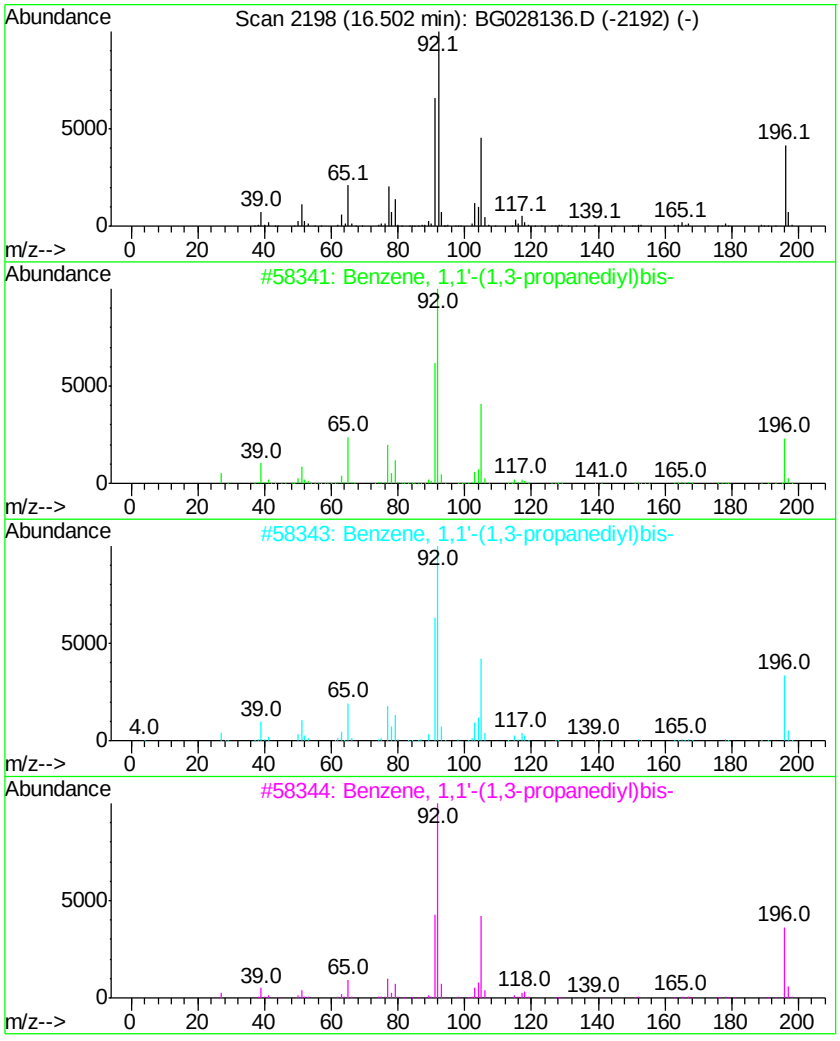
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

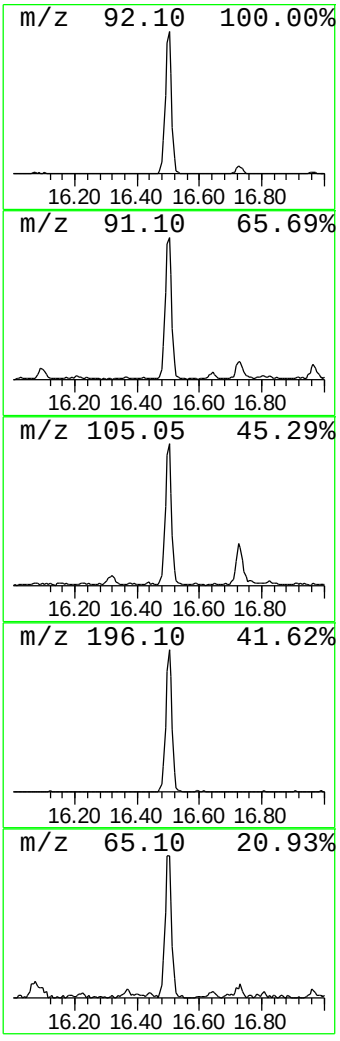
****

Peak Number 4 Benzene, 1,1'-(1,3-propaned... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.50	11.88 ng/ul	507762	Phenanthrene-d10	17.72

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,1'-(1,3-propanediyl)bis-	196	C15H16	001081-75-0	97
2	Benzene, 1,1'-(1,3-propanediyl)bis-	196	C15H16	001081-75-0	96
3	Benzene, 1,1'-(1,3-propanediyl)bis-	196	C15H16	001081-75-0	90
4	Benzene, 1,1'-(1,3-propanediyl)bis-	196	C15H16	001081-75-0	70
5	1-Chloro-6-phenylhexane	196	C12H17Cl	071434-68-9	64





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Sample : I4405-07
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : JJ2X4

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG080217.M
Quant Title : SVOA CALIBRATION

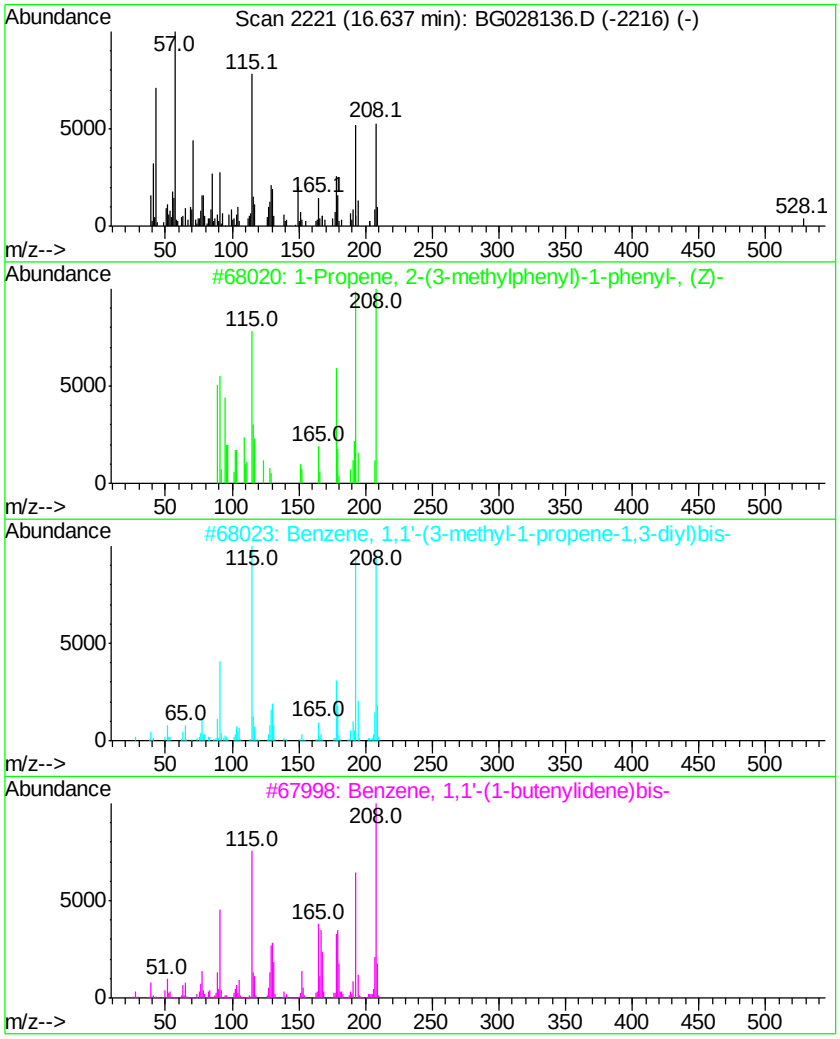
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

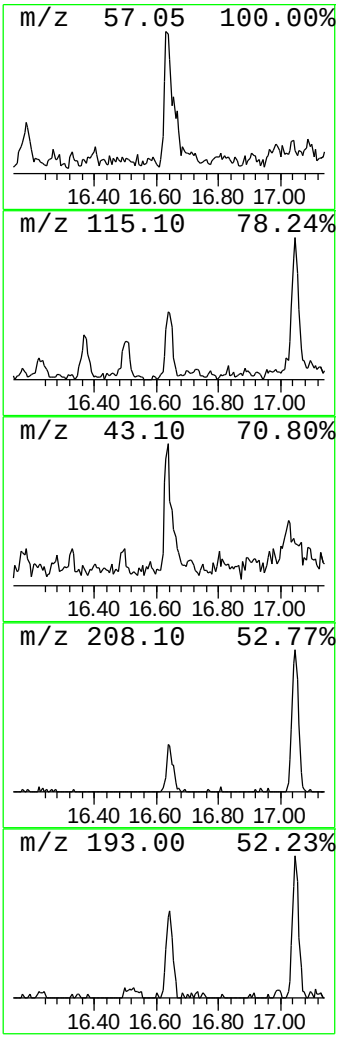
****

Peak Number 5 1-Propene, 2-(3-methylpheny... Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.64	2.51 ng/ul	107202	Phenanthrene-d10	17.72

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Propene, 2-(3-methylphenyl)-1-...	208	C16H16	1000138-72-6	70
2	Benzene, 1,1'-(3-methyl-1-propen...	208	C16H16	007614-93-9	64
3	Benzene, 1,1'-(1-butenylidene)bis-	208	C16H16	001726-14-3	53
4	1-Propene, 2-(2-methylphenyl)-1-...	208	C16H16	1000138-72-4	43
5	Benzene, 1,1'-(1,2-dimethyl-1,2-...	208	C16H16	000782-06-9	30





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ALS Vial : 35 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : JJ2X4

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG080217.M
Quant Title : SVOA CALIBRATION

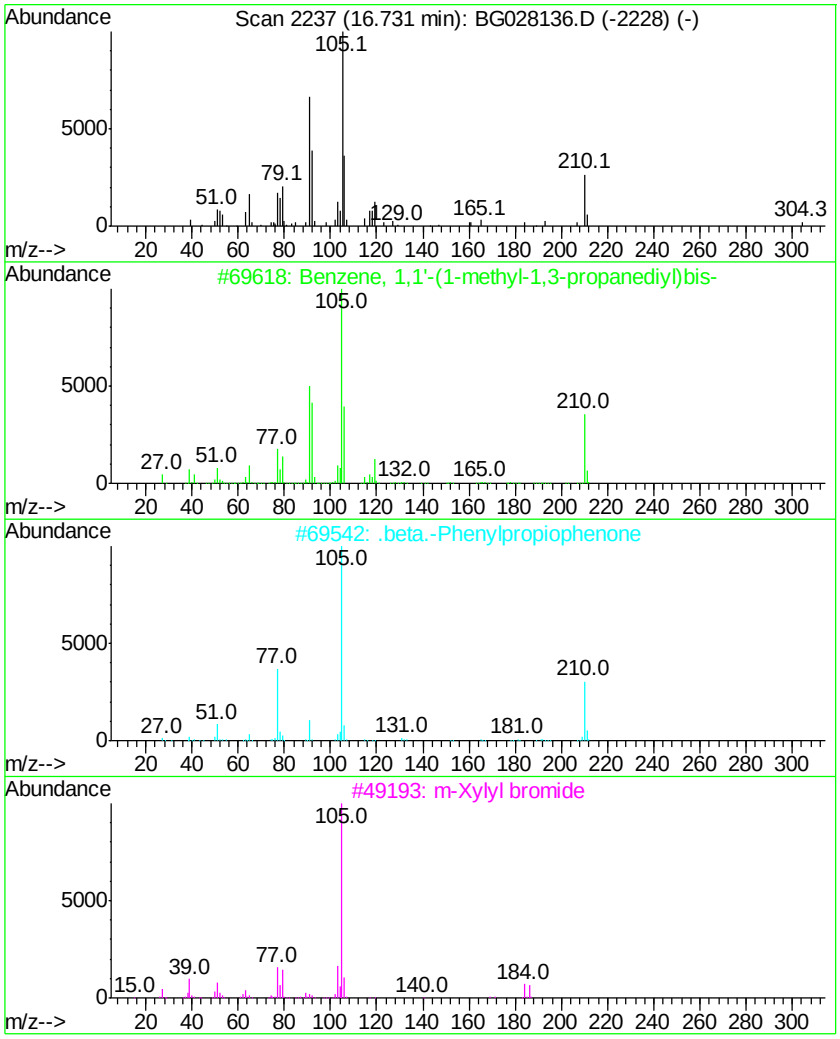
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

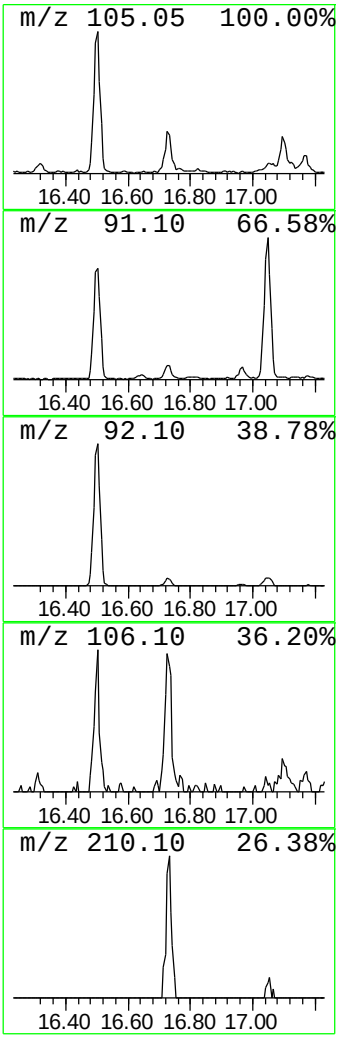
****

Peak Number 6 Benzene, 1,1'-(1-methyl-1,3... Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.73	2.43 ng/ul	104018	Phenanthrene-d10	17.72

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,1'-(1-methyl-1,3-prop...	210	C16H18	001520-44-1	91
2	.beta.-Phenylpropiophenone	210	C15H14O	001083-30-3	38
3	m-Xylyl bromide	184	C8H9Br	000620-13-3	35
4	Benzene, 1,1'-(1,4-butanediyl)bis-	210	C16H18	001083-56-3	30
5	Benzaldehyde, 3-ethoxy-4-(4-meth...	270	C17H18O3	351066-35-8	27





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Acq On : 3 Aug 2017 11:33
Operator : SJ/JU
Sample : I4405-07
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : JJ2X4

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG080217.M
Quant Title : SVOA CALIBRATION

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

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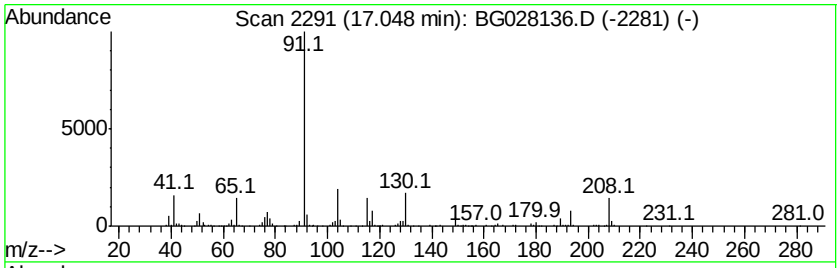
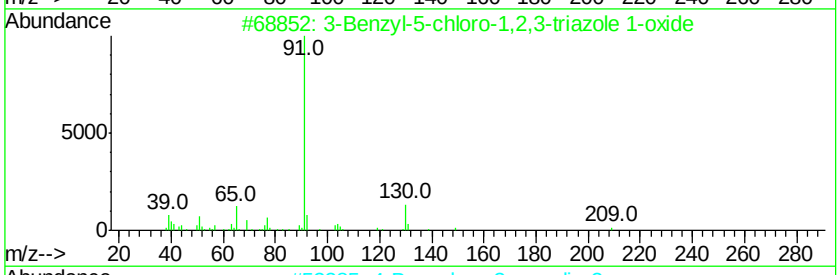
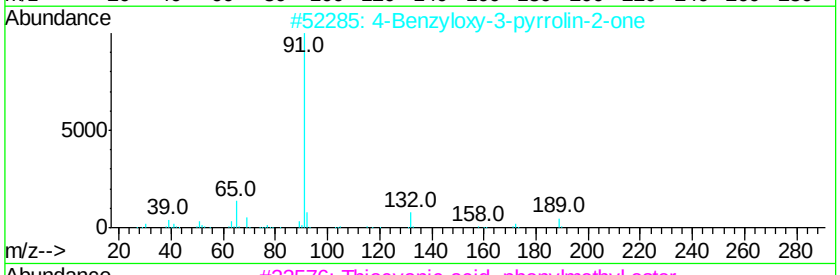
Peak Number 7 unknown-01 Concentration Rank 6

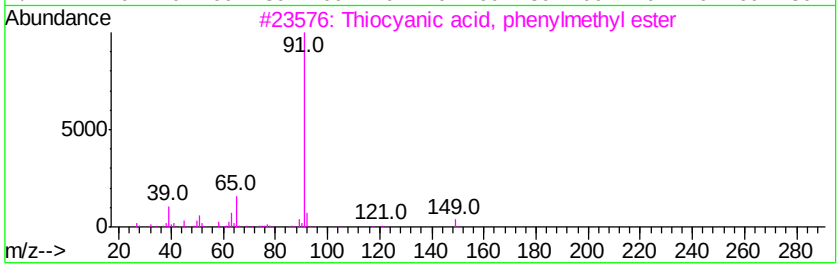
R.T.	EstConc	Area	Relative to ISTD	R.T.
17.05	14.40 ng/ul	615290	Phenanthrene-d10	17.72

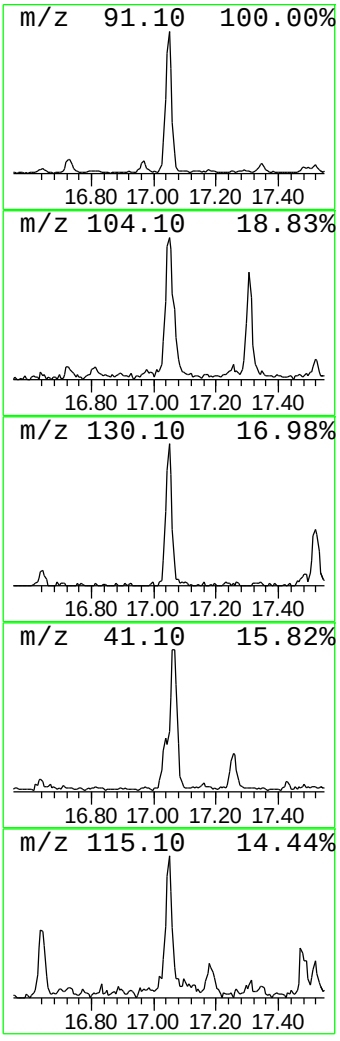
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Benzyl-5-chloro-1,2,3-triazole...	209	C9H8ClN3O	116932-67-3	41
2	4-Benzyloxy-3-pyrrolin-2-one	189	C11H11NO2	113896-95-0	25
3	Thiocyanic acid, phenylmethyl ester	149	C8H7NS	003012-37-1	25
4	Thiocyanic acid, phenylmethyl ester	149	C8H7NS	003012-37-1	22
5	N-Benzyl-1H-benzimidazole	208	C14H12N2	004981-92-4	22











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Data File : BG028136.D
Acq On : 3 Aug 2017 11:33
Operator : SJ/JU
Sample : I4405-07
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : JJ2X4

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG080217.M
Quant Title : SVOA CALIBRATION

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

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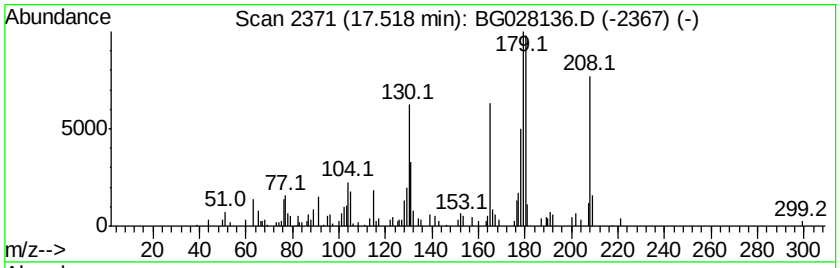
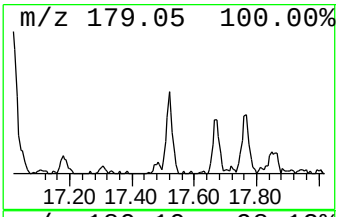
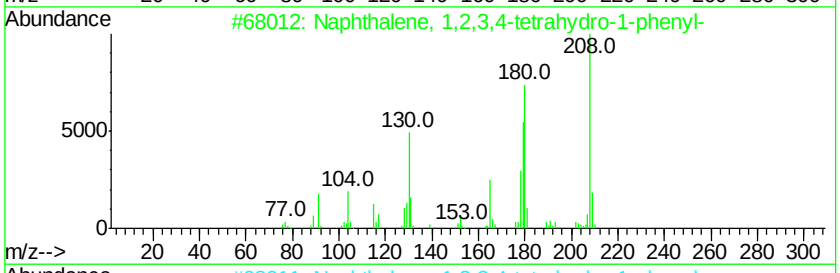
Peak Number 8 Naphthalene, 1,2,3,4-tetrah... Concentration Rank 22

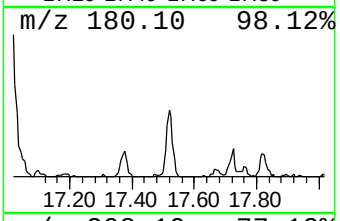
R.T.	EstConc	Area	Relative to ISTD	R.T.
17.52	2.68 ng/ul	114504	Phenanthrene-d10	17.72

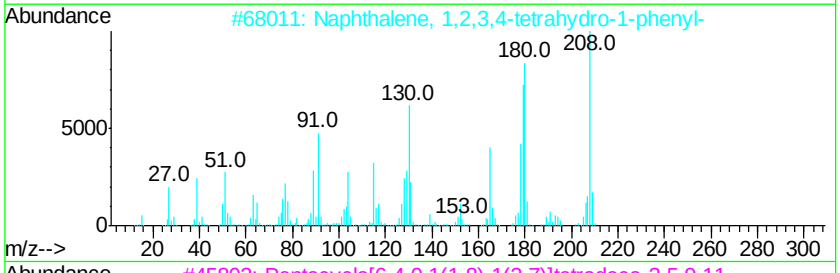
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,4-tetrahydro-...	208	C16H16	003018-20-0	90
2	Naphthalene, 1,2,3,4-tetrahydro-...	208	C16H16	003018-20-0	90
3	Pentacyclo[6.4.0.1(1,8).1(2,7)]t...	180	C14H12	086120-84-5	83
4	Naphthalene, 1,2,3,4-tetrahydro-...	208	C16H16	003018-20-0	70
5	Dibenzosuberone	208	C15H12O	001210-35-1	58

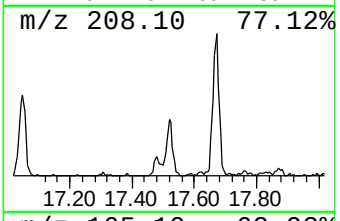


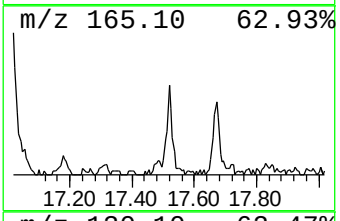


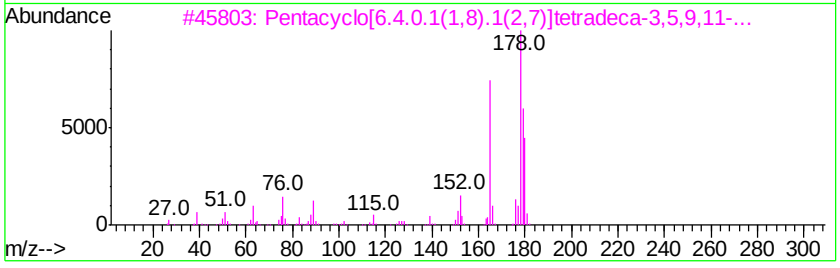


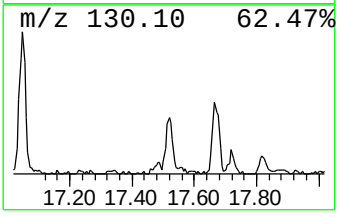












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Data File : BG028136.D
Acq On : 3 Aug 2017 11:33
Operator : SJ/JU
Sample : I4405-07
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : JJ2X4

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG080217.M
Quant Title : SVOA CALIBRATION

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

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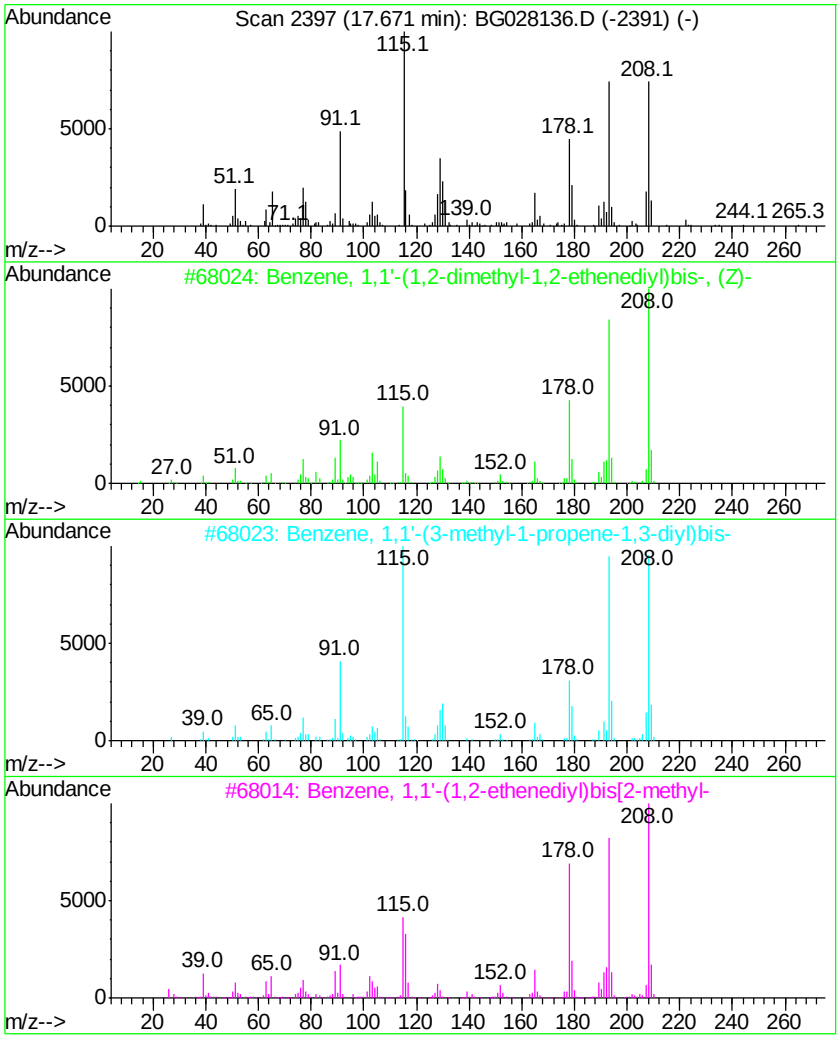
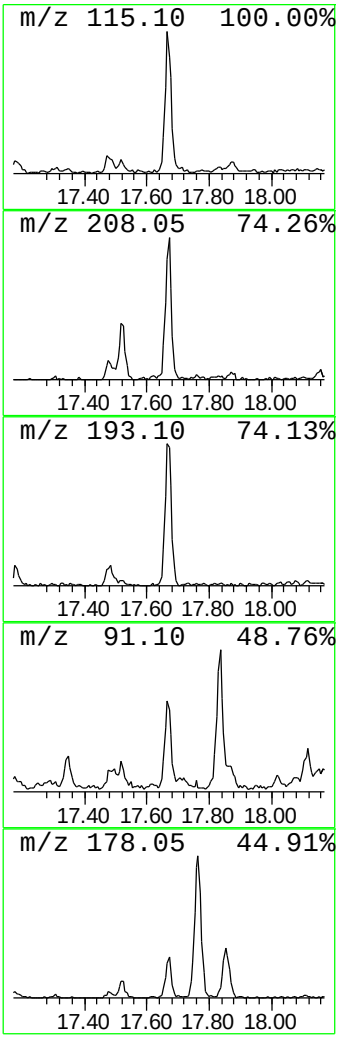
Peak Number 9 Benzene, 1,1'-(1,2-dimethyl... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.67	6.67 ng/ul	285047	Phenanthrene-d10	17.72

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,1'-(1,2-dimethyl-1,2-...	208	C16H16	000782-05-8	95
2	Benzene, 1,1'-(3-methyl-1-propen...	208	C16H16	007614-93-9	90
3	Benzene, 1,1'-(1,2-ethenediyl)bi...	208	C16H16	010311-74-7	86
4	Benzene, 1,1'-(1,2-dimethyl-1,2-...	208	C16H16	000782-06-9	60
5	1-Propene, 2-(2-methylphenyl)-1-...	208	C16H16	1000138-72-4	60





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Acq On : 3 Aug 2017 11:33
Operator : SJ/JU
Sample : I4405-07
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : JJ2X4

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG080217.M
Quant Title : SVOA CALIBRATION

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

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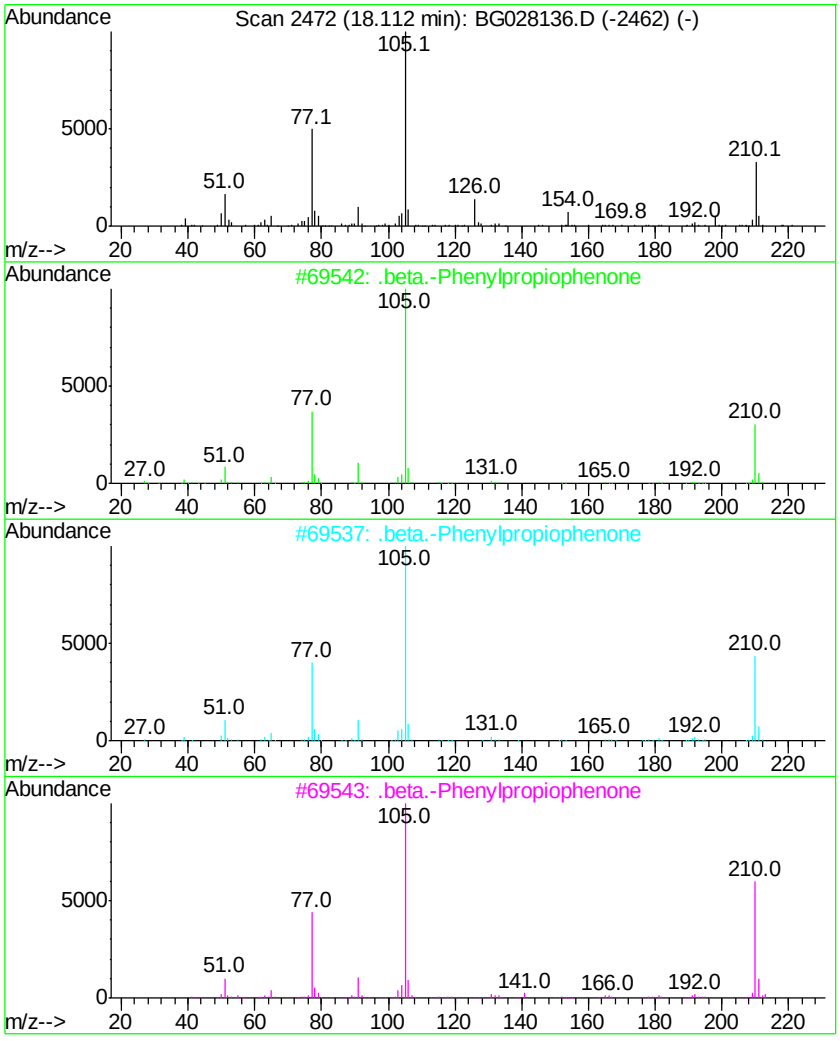
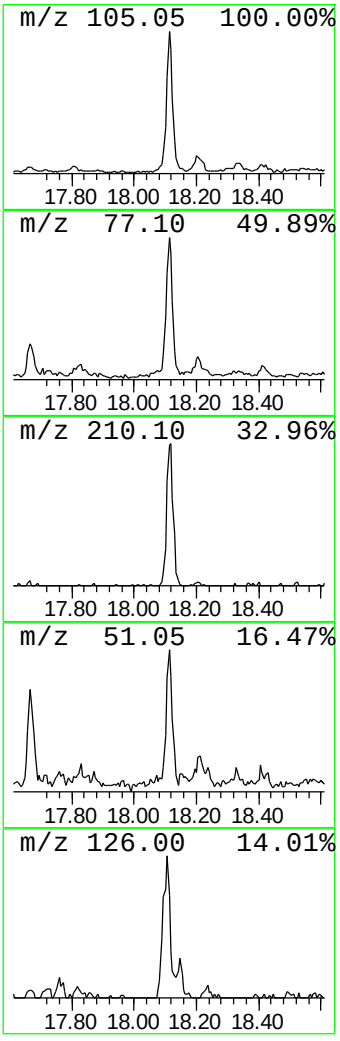
Peak Number 10 .beta.-Phenylpropiophenone Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.11	6.84 ng/ul	292398	Phenanthrene-d10	17.72

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	.beta.-Phenylpropiophenone	210	C15H14O	001083-30-3	94
2	.beta.-Phenylpropiophenone	210	C15H14O	001083-30-3	93
3	.beta.-Phenylpropiophenone	210	C15H14O	001083-30-3	81
4	N-Benzylbenzamide	211	C14H13NO	001485-70-7	53
5	Benzoic acid, phenyl ester	198	C13H10O2	000093-99-2	49





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ALS Vial : 35 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : JJ2X4

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG080217.M
Quant Title : SVOA CALIBRATION

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

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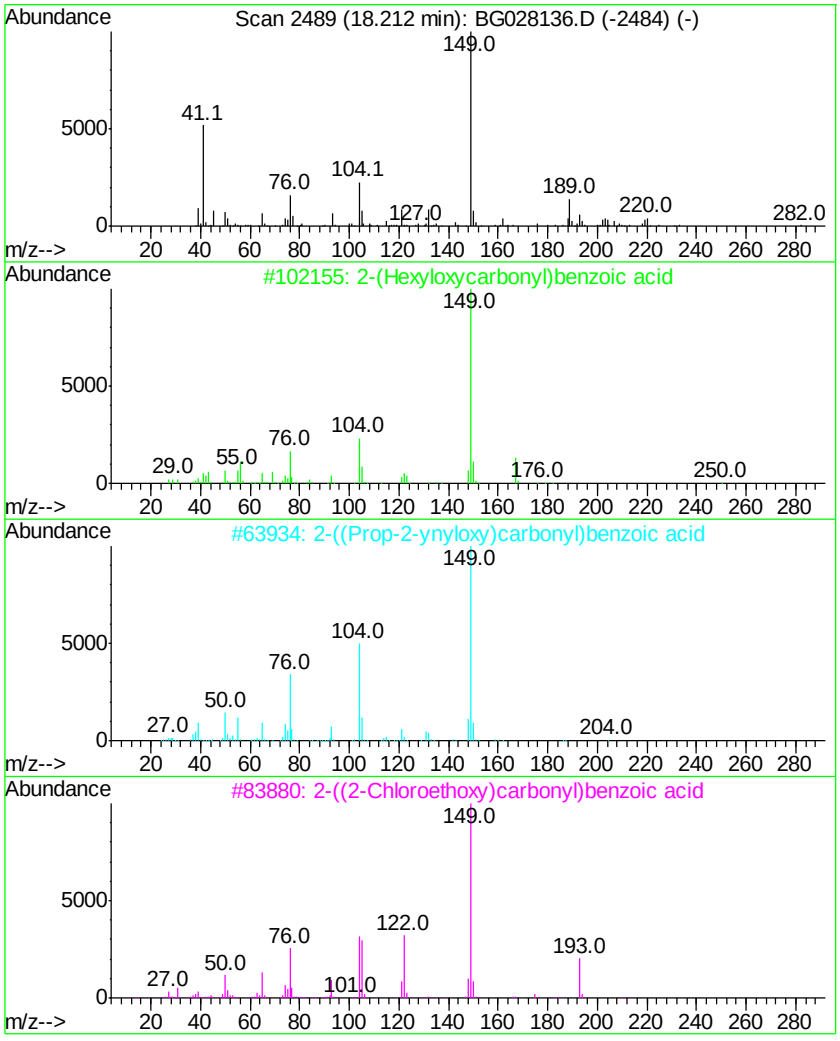
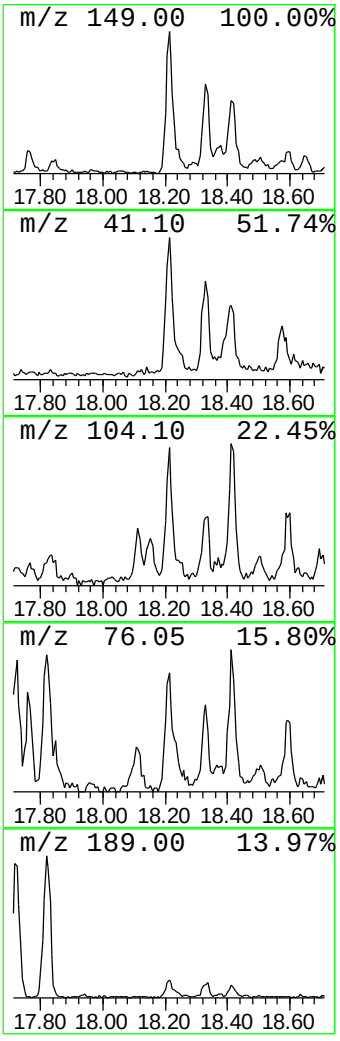
Peak Number 11 2-(Hexyloxycarbonyl)benzoic... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.21	7.11 ng/ul	303745	Phenanthrene-d10	17.72

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-(Hexyloxycarbonyl)benzoic acid	250	C14H18O4	1000373-63-0	59
2	2-((Prop-2-ynyloxy)carbonyl)benz...	204	C11H8O4	1000373-53-6	59
3	2-((2-Chloroethoxy)carbonyl)benz...	228	C10H9ClO4	1000373-53-5	59
4	2-((But-3-enyloxy)carbonyl)benzo...	220	C12H12O4	1000373-53-4	59
5	2-(Octyloxycarbonyl)benzoic acid	278	C16H22O4	1000373-53-3	53





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Sample : I4405-07
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : JJ2X4

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Quant Title : SVOA CALIBRATION

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

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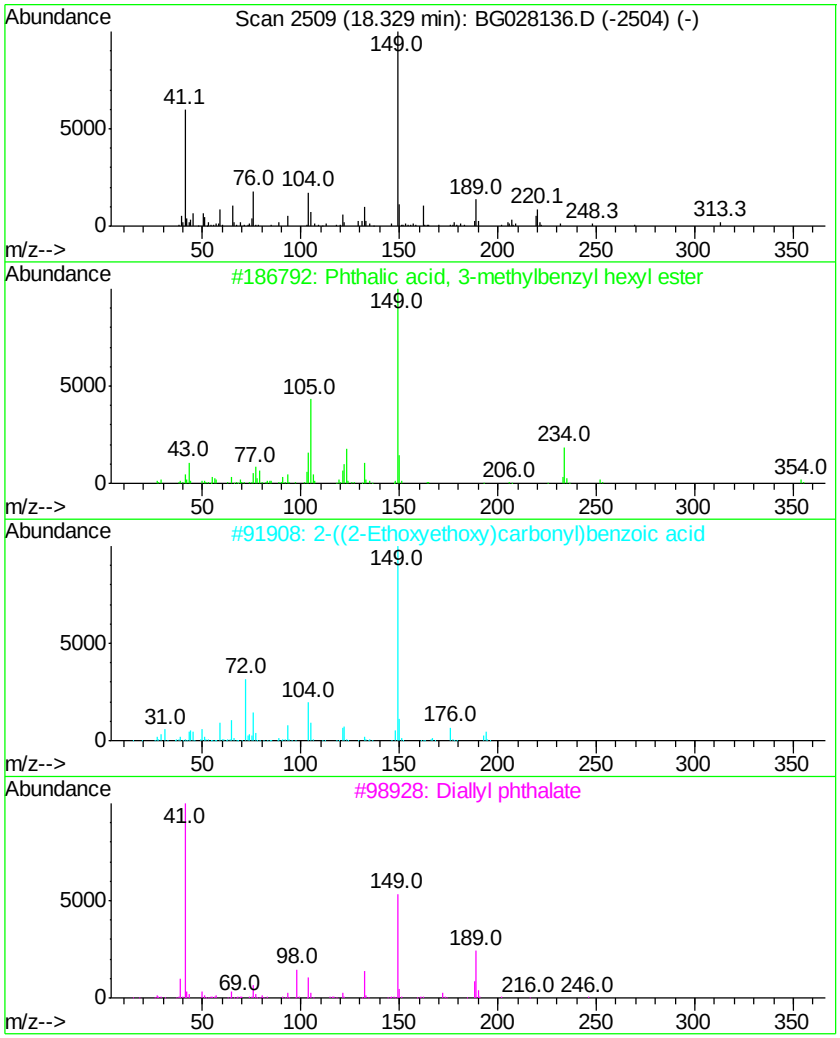
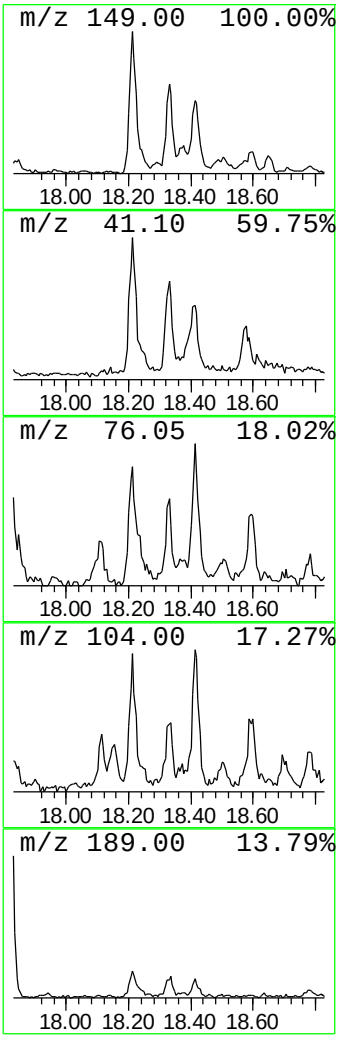
Peak Number 12 Phthalic acid, 3-methylbenz... Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.33	2.01 ng/ul	86095	Phenanthrene-d10	17.72

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phthalic acid, 3-methylbenzyl he...	354	C22H26O4	1000371-10-4	59
2	2-((2-Ethoxyethoxy)carbonyl)benz...	238	C12H14O5	1000373-82-8	53
3	Diallyl phthalate	246	C14H14O4	000131-17-9	53
4	Diallyl phthalate	246	C14H14O4	000131-17-9	53
5	Phthalic acid, oct-3-yl pentyl e...	348	C21H32O4	1000377-71-4	53





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Sample : I4405-07
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : JJ2X4

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG080217.M
Quant Title : SVOA CALIBRATION

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

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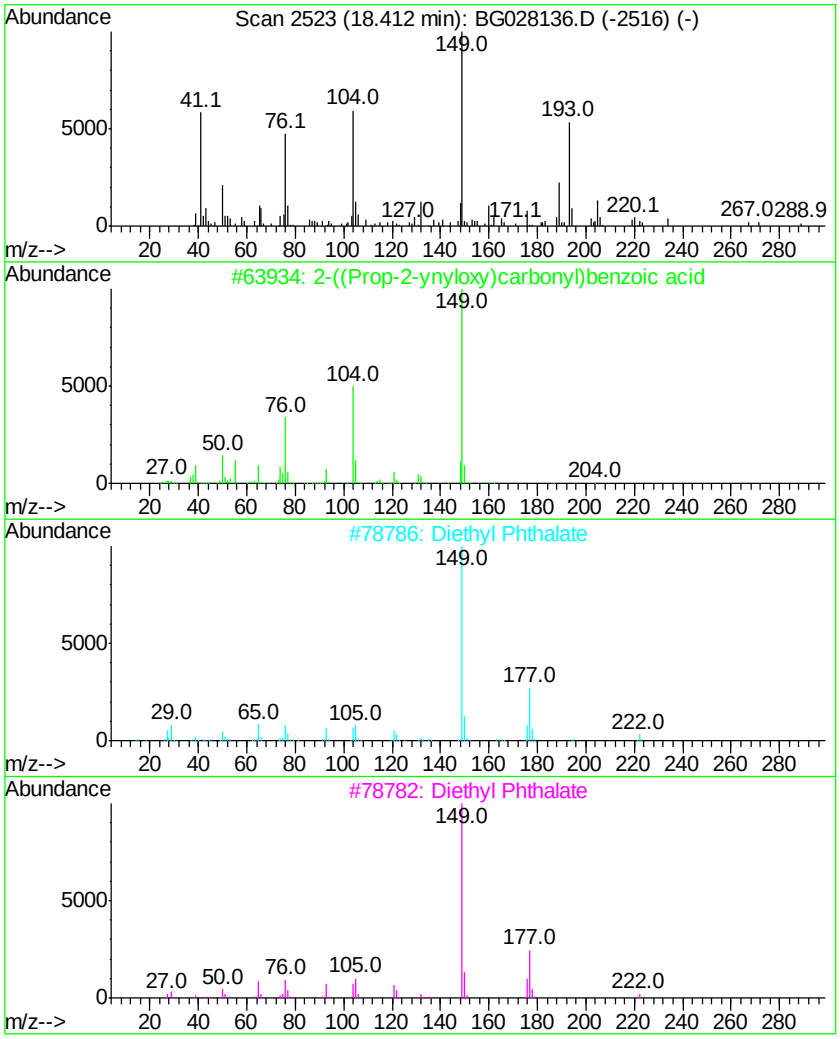
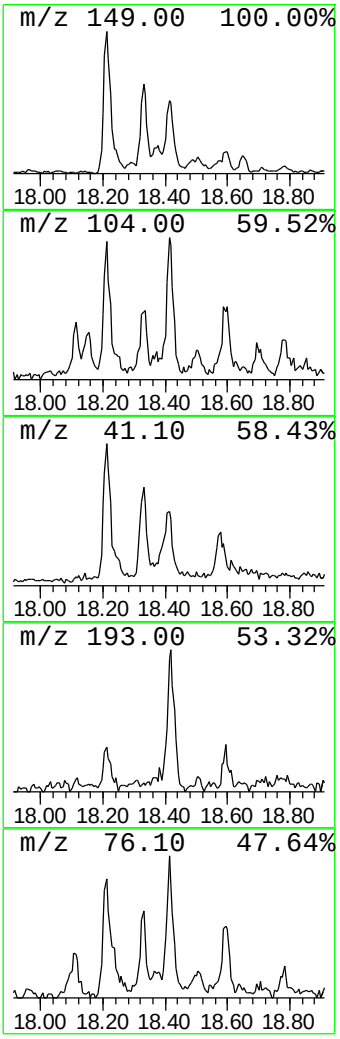
Peak Number 13 unknown-02 Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.41	5.11 ng/ul	218556	Phenanthrene-d10	17.72

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-((Prop-2-ynyloxy)carbonyl)benz...	204	C11H8O4	1000373-53-6	47
2	Diethyl Phthalate	222	C12H14O4	000084-66-2	22
3	Diethyl Phthalate	222	C12H14O4	000084-66-2	22
4	Phthalic acid, butyl 3,3-dimethy...	306	C18H26O4	1000357-00-2	16
5	1H-S-Triazolo[1,5-a]pyridin-4-iu...	149	C7H7N3O	013980-64-8	14





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Acq On : 3 Aug 2017 11:33
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Sample : I4405-07
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : JJ2X4

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Quant Title : SVOA CALIBRATION

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

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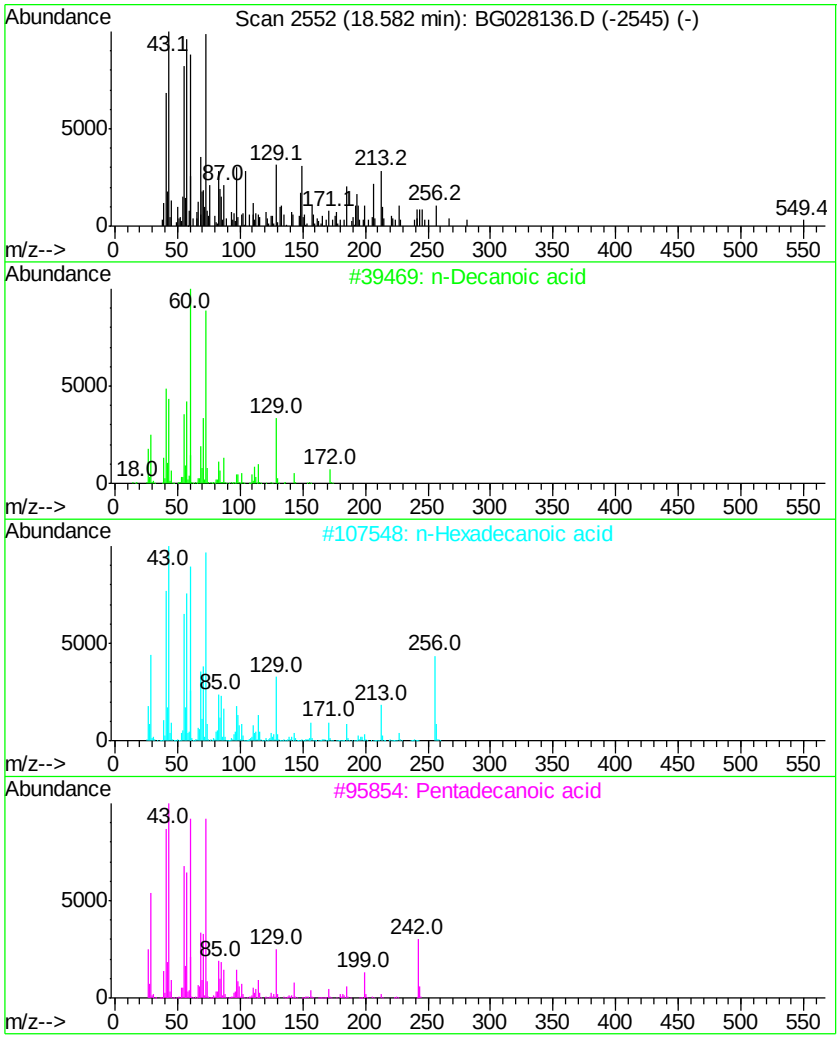
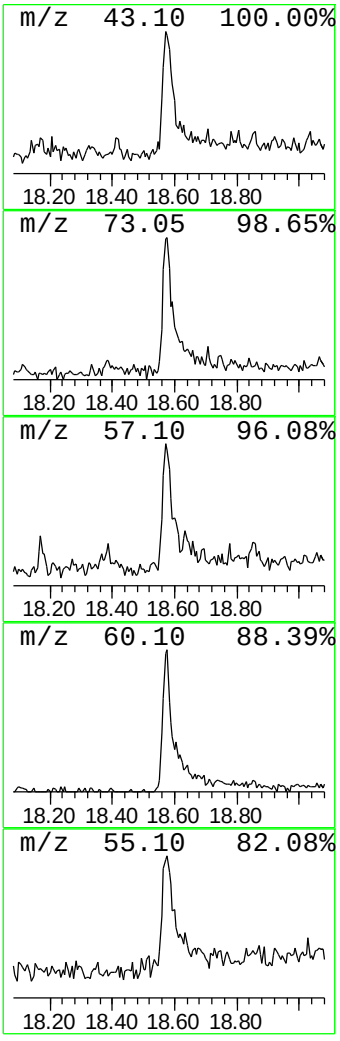
Peak Number 14 n-Decanoic acid Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.58	5.89 ng/ul	251948	Phenanthrene-d10	17.72

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Decanoic acid	172	C10H20O2	000334-48-5	91
2	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	64
3	Pentadecanoic acid	242	C15H30O2	001002-84-2	58
4	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	58
5	Tridecanoic acid	214	C13H26O2	000638-53-9	53





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Sample : I4405-07
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ALS Vial : 35 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : JJ2X4

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Quant Title : SVOA CALIBRATION

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

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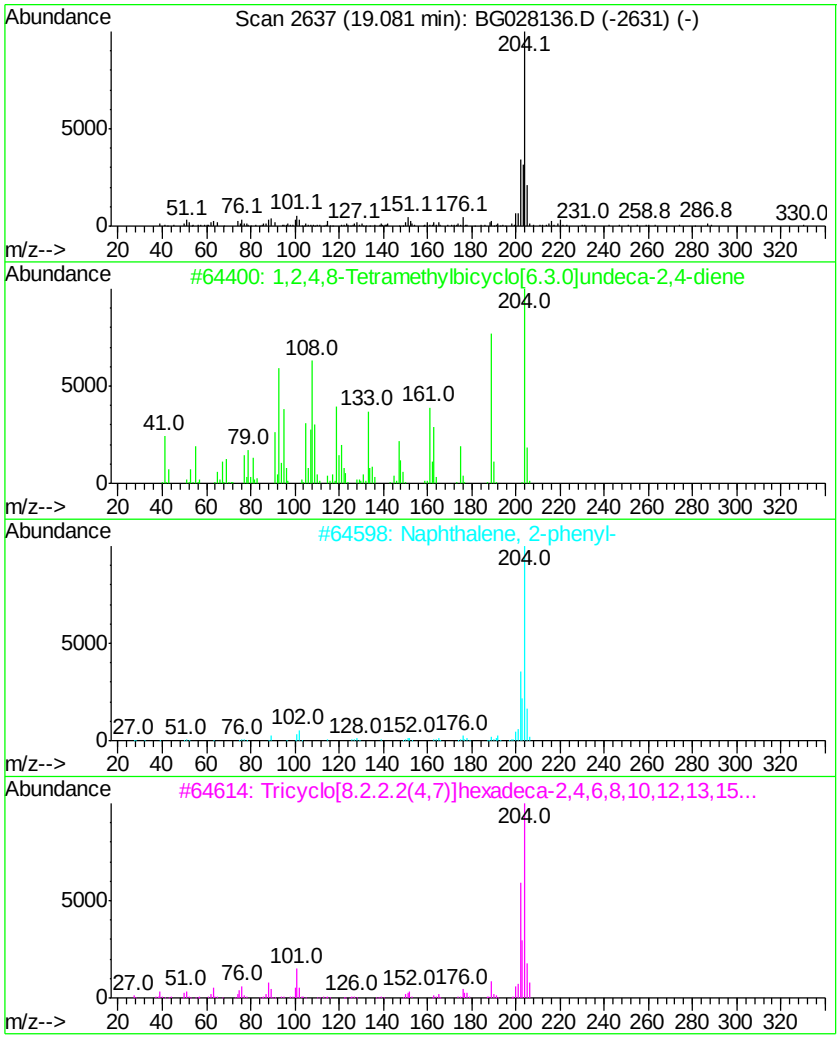
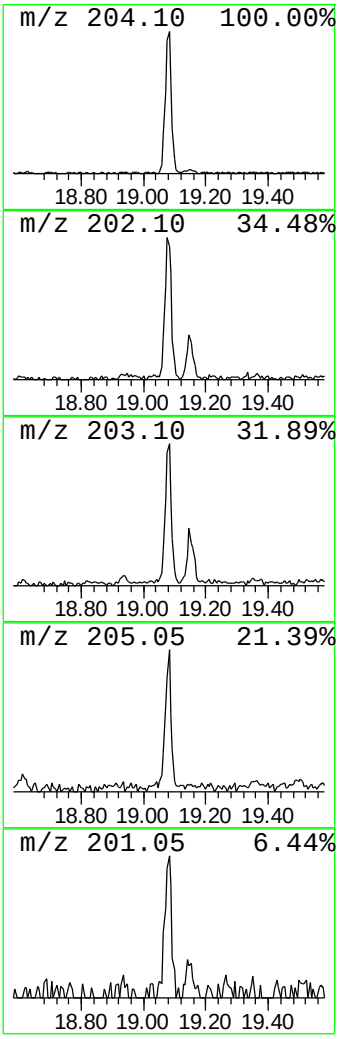
Peak Number 15 1,2,4,8-Tetramethylbicyclo[... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.08	4.27 ng/ul	182675	Phenanthrene-d10	17.72

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,2,4,8-Tetramethylbicyclo[6.3.0...	204	C15H24	137235-51-9	92
2	Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	81
3	Tricyclo[8.2.2.2(4,7)]hexadeca-2...	204	C16H12	006572-60-7	81
4	5,16[1',2']:8,13[1'',2'']-Dibenz...	408	C32H24	005672-97-9	80
5	Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	76





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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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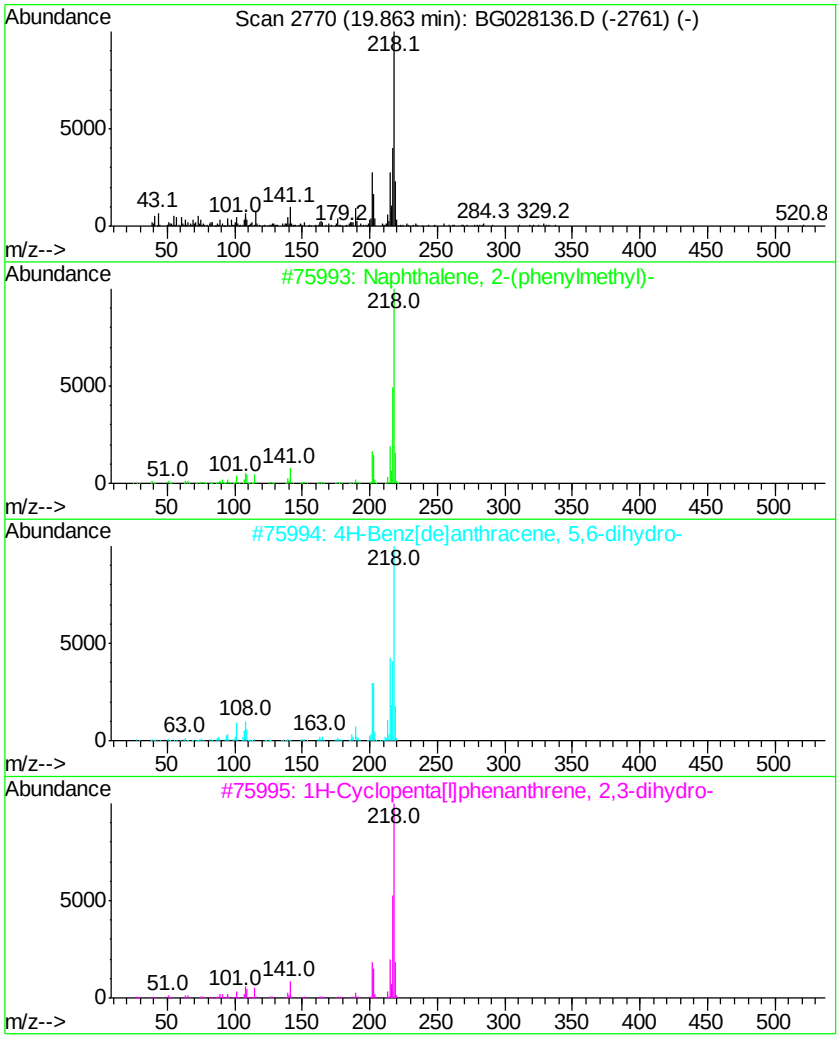
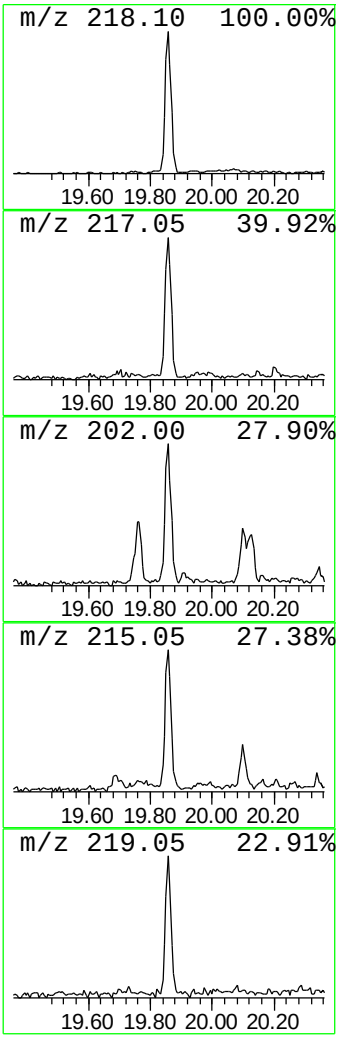
Peak Number 16 Naphthalene, 2-(phenylmethyl)- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.86	9.64 ng/ul	412107	Phenanthrene-d10	17.72

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2-(phenylmethyl)-	218	C17H14	000613-59-2	83
2	4H-Benz[de]anthracene, 5,6-dihydro-	218	C17H14	004389-09-7	76
3	1H-Cyclopenta[l]phenanthrene, 2,...	218	C17H14	000723-98-8	70
4	6-Methyl-8-phenylbenzocyclohepte...	246	C18H14O	1000211-20-7	56
5	8-Amino-2,6-dimethoxylepidine	218	C12H14N2O2	074509-65-2	53





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ALS Vial : 35 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : JJ2X4

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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
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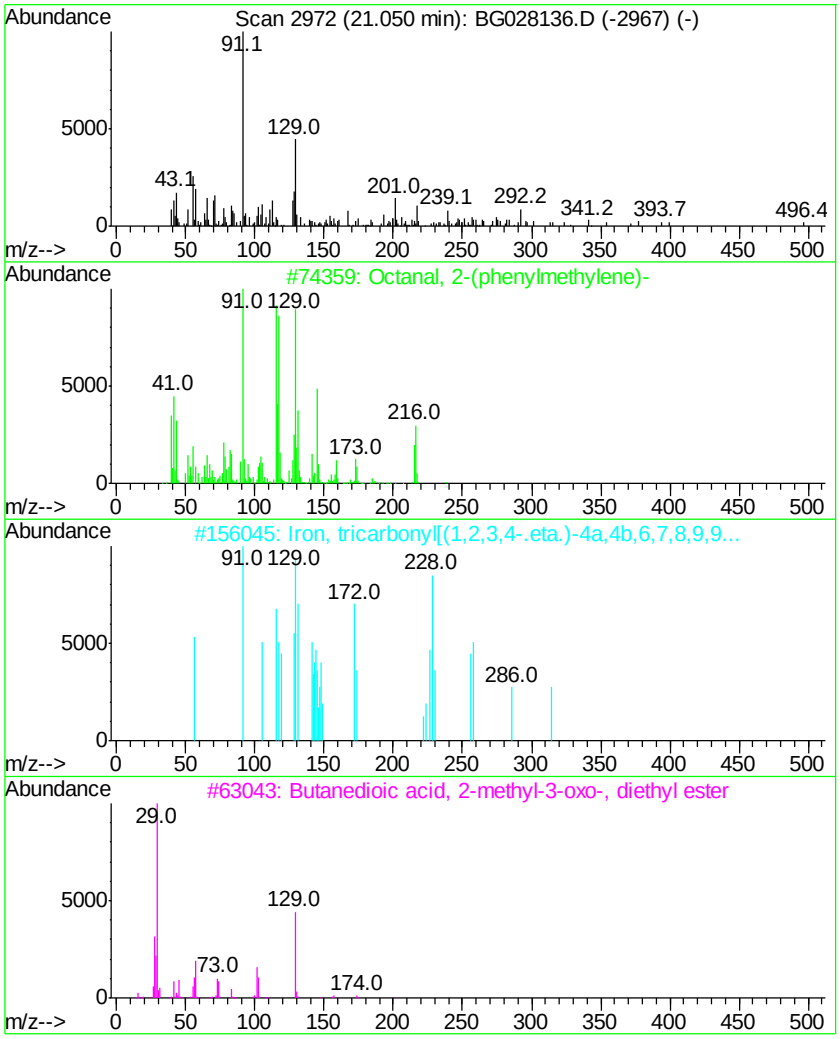
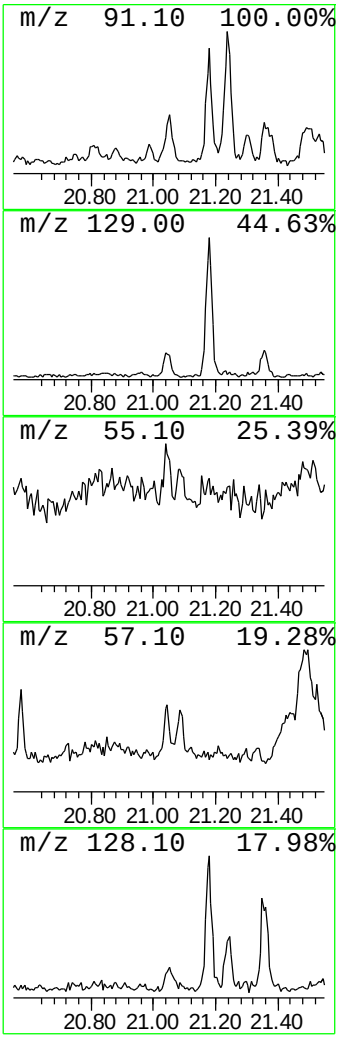
Peak Number 17 unknown-03 Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.05	2.16 ng/ul	108671	Chrysene-d12	22.05

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octanal, 2-(phenylmethylene)-	216	C15H20O	000101-86-0	38
2	Iron, tricarbonyl[(1,2,3,4-.eta....	314	C16H18FeO3	056784-34-0	35
3	Butanedioic acid, 2-methyl-3-oxo...	202	C9H14O5	000759-65-9	17
4	1-Methoxy-5-phenyl-1-trimethysil...	248	C15H24OSi	1000196-36-6	9
5	5,9-Methanobenzocycloocten-11-on...	307	C19H17NO3	101401-33-6	9





Data Path : Z:\HPCHEM1\BNA G\DATA\BG080217\
Data File : BG028136.D
Acq On : 3 Aug 2017 11:33
Operator : SJ/JU
Sample : I4405-07
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : JJ2X4

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG080217.M
Quant Title : SVOA CALIBRATION

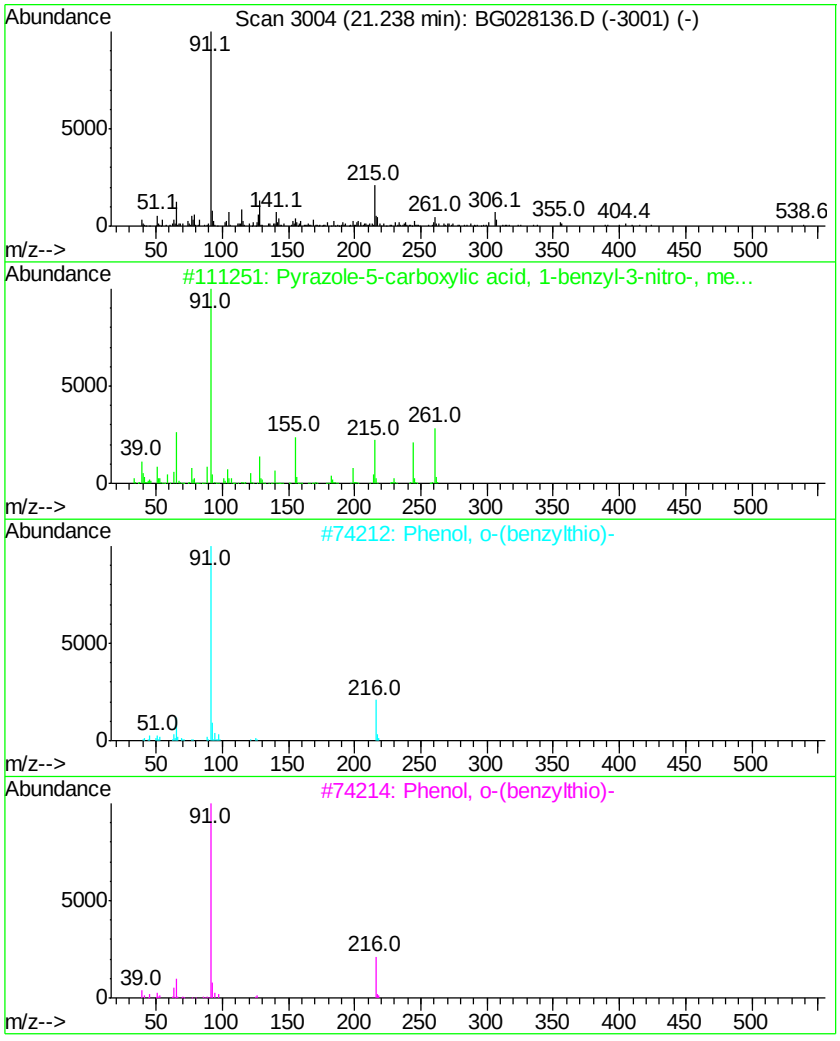
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

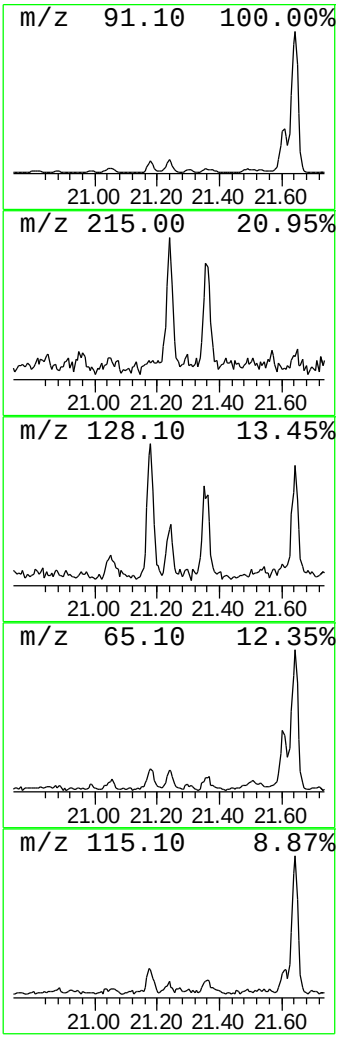
****

Peak Number 19 unknown-04 Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.24	2.85 ng/ul	143696	Chrysene-d12	22.05

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pyrazole-5-carboxylic acid, 1-be...	261	C12H11N3O4	1000271-76-6	36
2	Phenol, o-(benzylthio)-	216	C13H12OS	024312-63-8	14
3	Phenol, o-(benzylthio)-	216	C13H12OS	024312-63-8	14
4	Bibenzyl, 3-chloro-	216	C14H13Cl	034176-92-6	14
5	N .epsilon.-Carbobenzyloxy-L-lysine	280	C14H20N2O4	001155-64-2	12





Library Search Compound Report

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Data File : BG028136.D
Acq On : 3 Aug 2017 11:33
Operator : SJ/JU
Sample : I4405-07
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : JJ2X4

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG080217.M
Quant Title : SVOA CALIBRATION

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

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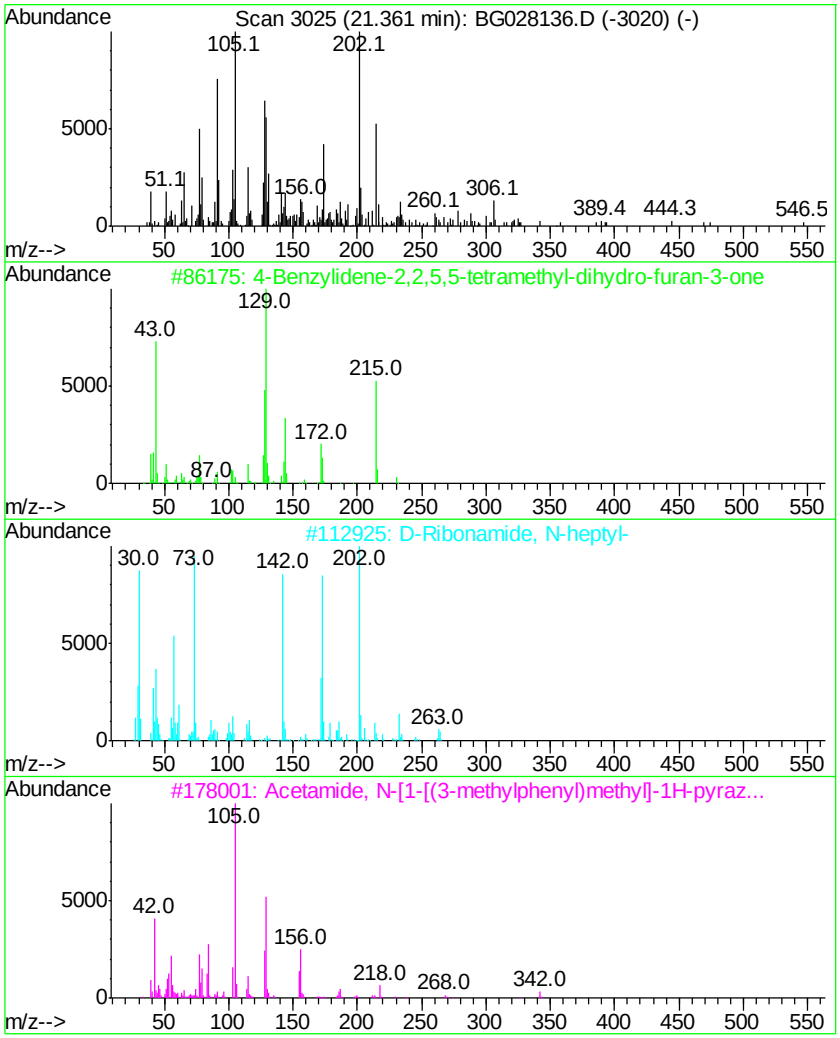
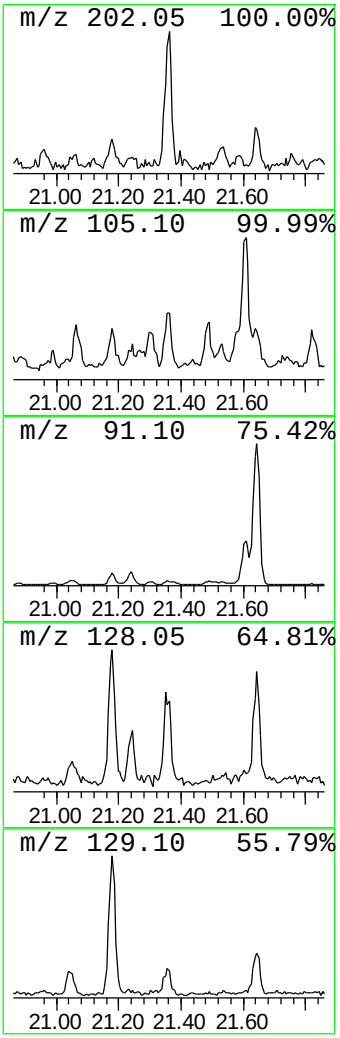
Peak Number 20 unknown-05 Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.36	3.78 ng/ul	190319	Chrysene-d12	22.05

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-Benzylidene-2,2,5,5-tetramethy...	230	C15H18O2	1000301-02-8	27
2	D-Ribonamide, N-heptyl-	263	C12H25NO5	1000158-56-2	20
3	Acetamide, N-[1-[(3-methylphenyl...	342	C16H18N6OS	1000319-53-9	18
4	Benzene, 1,1'-(1,3-butadiyne-1,4...	202	C16H10	000886-66-8	18
5	1,2,3,6-Tetrahydro-oxypyridine, ...	203	C13H17NO	095969-41-8	15





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Data File : BG028136.D
Acq On : 3 Aug 2017 11:33
Operator : SJ/JU
Sample : I4405-07
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : JJ2X4

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG080217.M
Quant Title : SVOA CALIBRATION

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

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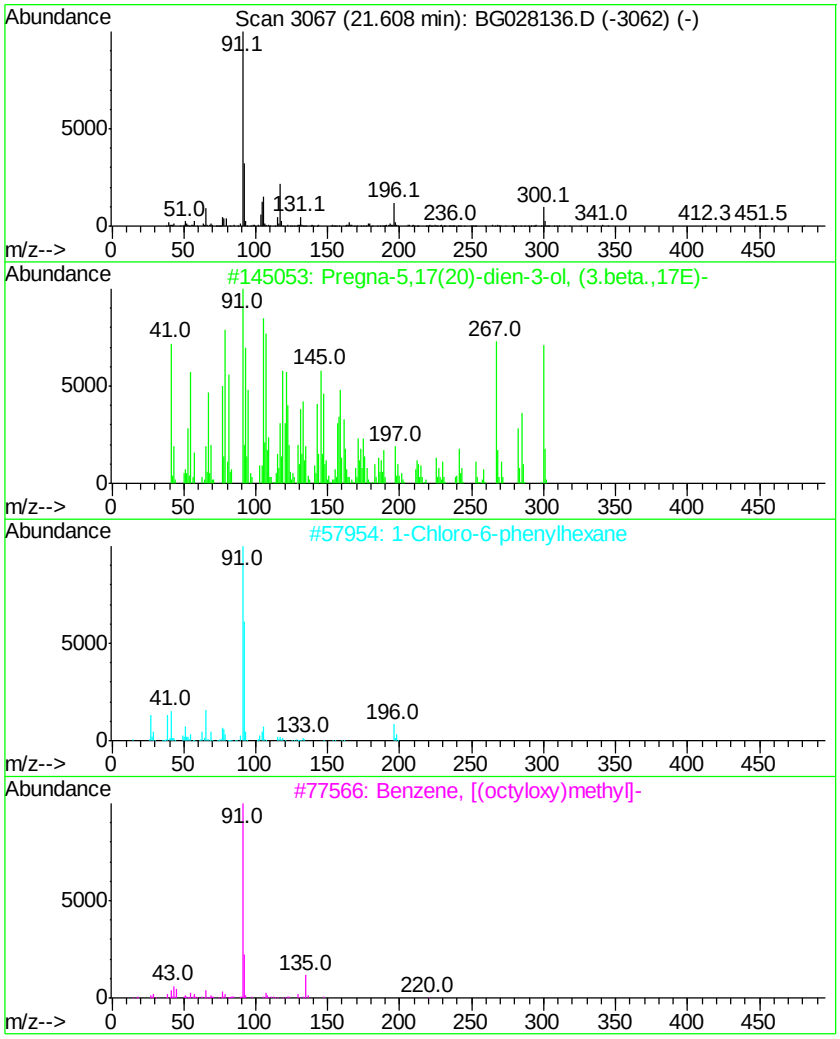
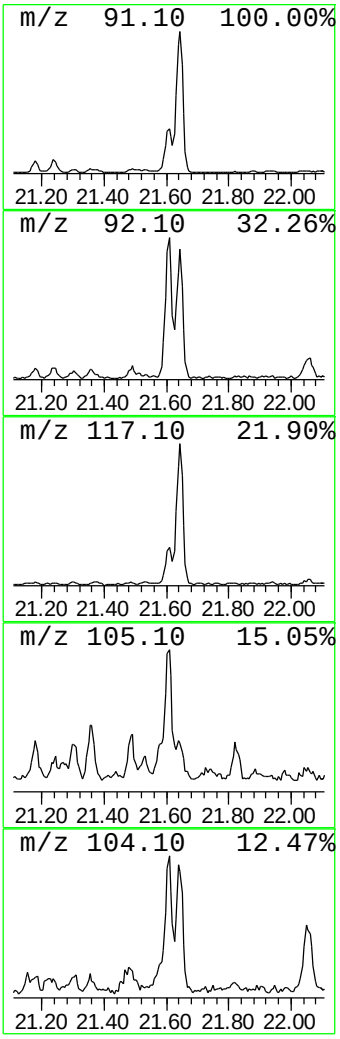
Peak Number 21 Pregna-5,17(20)-dien-3-ol, ... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.61	8.12 ng/ul	408817	Chrysene-d12	22.05

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pregna-5,17(20)-dien-3-ol, (3.be...	300	C21H32O	001159-25-7	59
2	1-Chloro-6-phenylhexane	196	C12H17Cl	071434-68-9	25
3	Benzene, [(octyloxy)methyl]-	220	C15H24O	054852-64-1	22
4	Androstan-17-one, 3-ethyl-3-hydr...	318	C21H34O2	057344-99-7	22
5	(3-Bromo-1-methylpropoxymethyl)b...	242	C11H15BrO	051666-29-6	22





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Data File : BG028136.D
Acq On : 3 Aug 2017 11:33
Operator : SJ/JU
Sample : I4405-07
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : JJ2X4

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG080217.M
Quant Title : SVOA CALIBRATION

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

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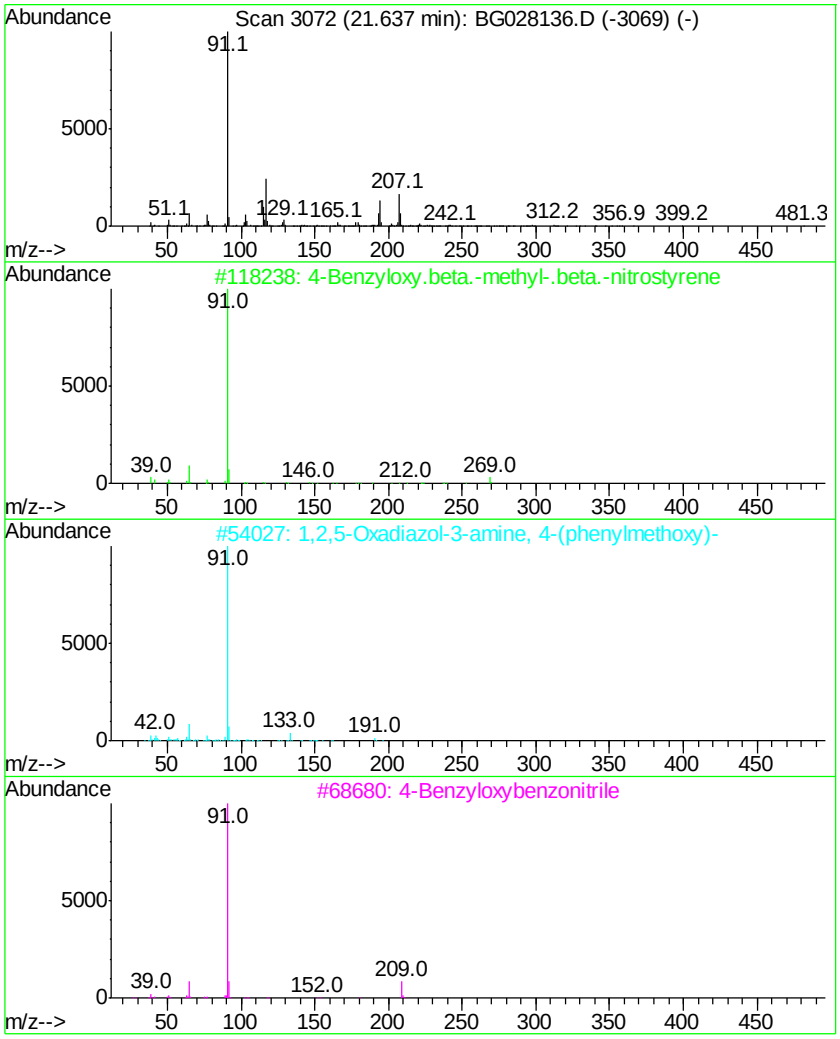
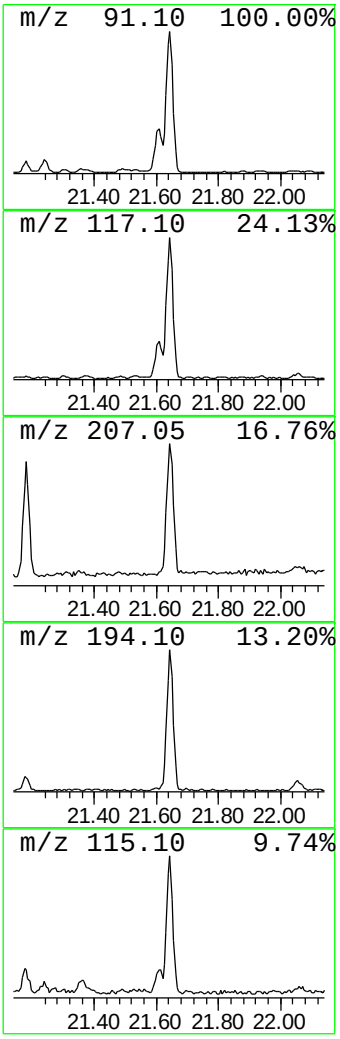
Peak Number 22 unknown-06 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.64	26.44 ng/ul	1331000	Chrysene-d12	22.05

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-Benzyloxy.beta.-methyl-.beta.-...	269	C16H15NO3	007176-38-7	10
2	1,2,5-Oxadiazol-3-amine, 4-(phen...	191	C9H9N3O2	1000337-58-8	10
3	4-Benzyloxybenzonitrile	209	C14H11NO	052805-36-4	10
4	2,3-Butanediol, 1,4-bis[(benzylo...	388	C20H24N2O6	1000163-34-1	10
5	Alpha-phenyl-alpha-tropylacetald...	378	C22H22N2O2S	022532-16-7	10





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Data File : BG028136.D
Acq On : 3 Aug 2017 11:33
Operator : SJ/JU
Sample : I4405-07
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : JJ2X4

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG080217.M
Quant Title : SVOA CALIBRATION

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

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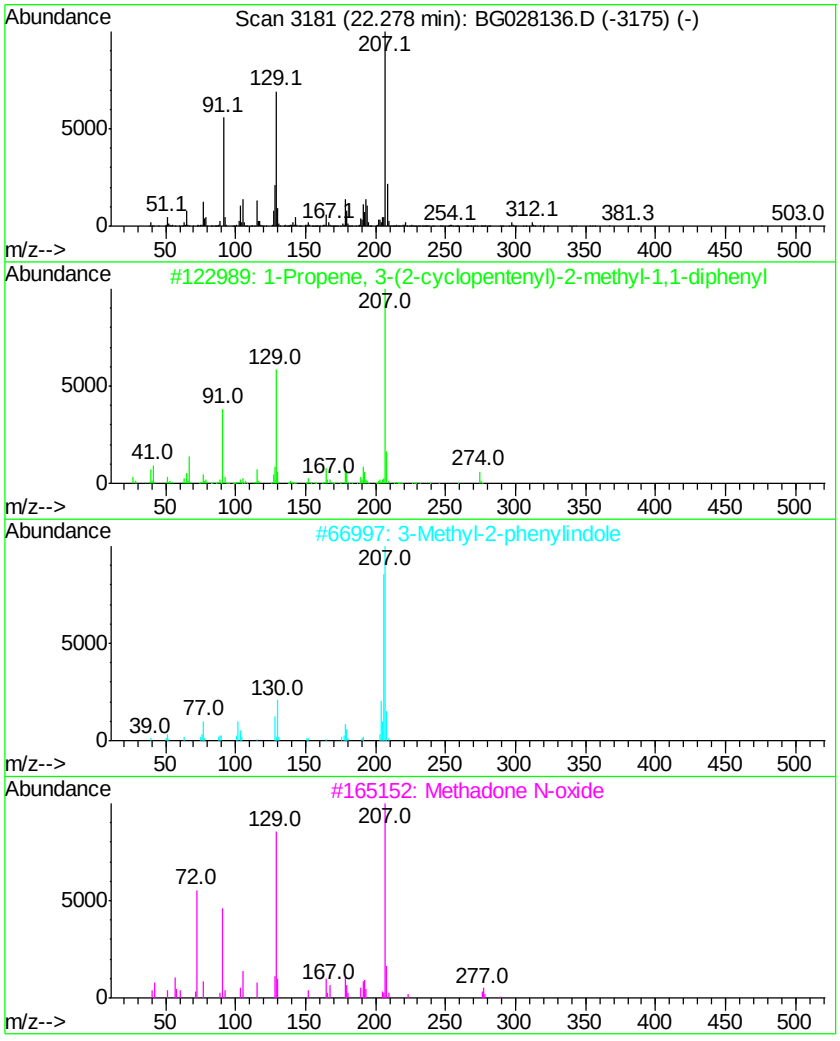
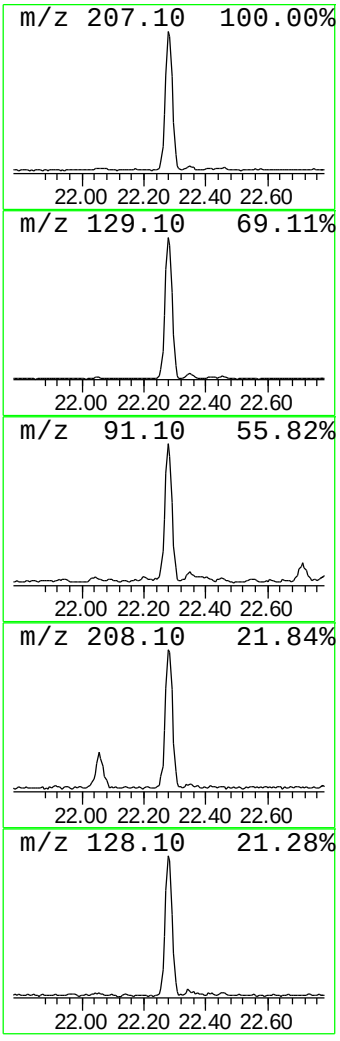
Peak Number 23 1-Propene, 3-(2-cyclopenten... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.28	33.62 ng/ul	1692720	Chrysene-d12	22.05

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Propene, 3-(2-cyclopentenyl)-2...	274	C21H22	1000154-23-3	83
2	3-Methyl-2-phenylindole	207	C15H13N	010257-92-8	46
3	Methadone N-oxide	325	C21H27NO2	1000120-80-7	43
4	9,10-Methanoanthracen-11-ol, 9,1...	250	C18H18O	126615-74-5	38
5	Hexahydropyridine, 1-methyl-4-[4...	207	C12H17NO2	094427-47-1	38





Data Path : Z:\HPCHEM1\BNA G\DATA\BG080217\
Data File : BG028136.D
Acq On : 3 Aug 2017 11:33
Operator : SJ/JU
Sample : I4405-07
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : JJ2X4

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG080217.M
Quant Title : SVOA CALIBRATION

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

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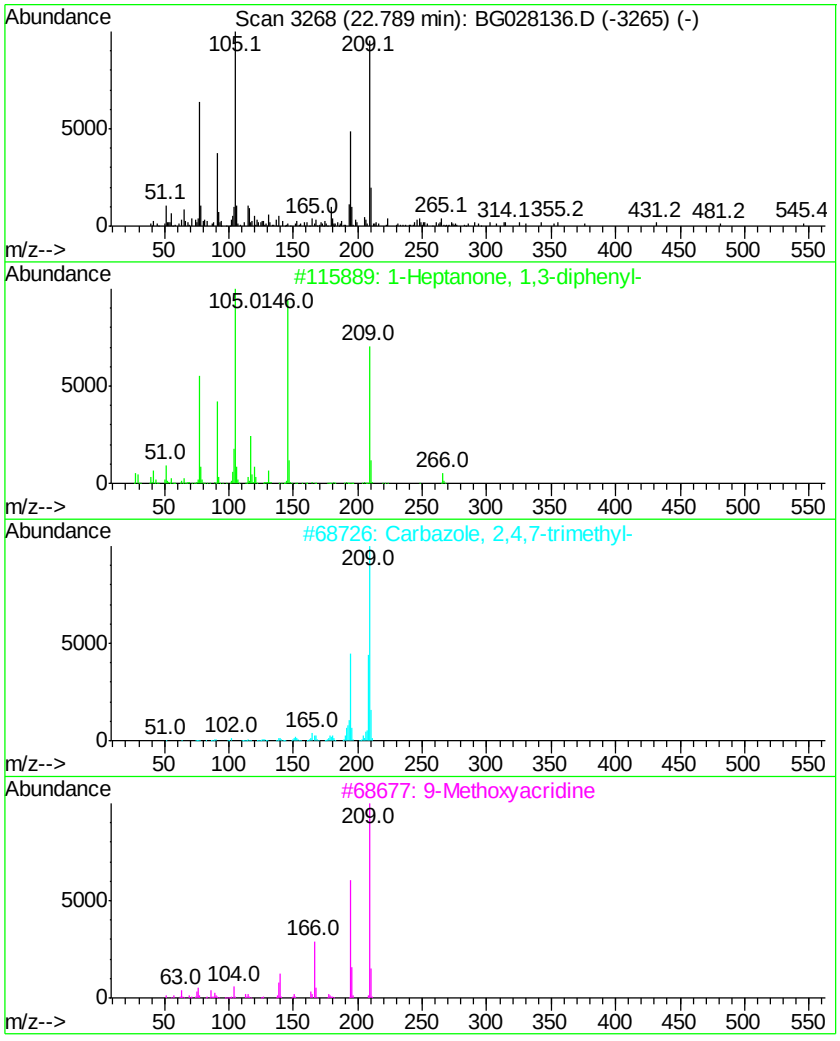
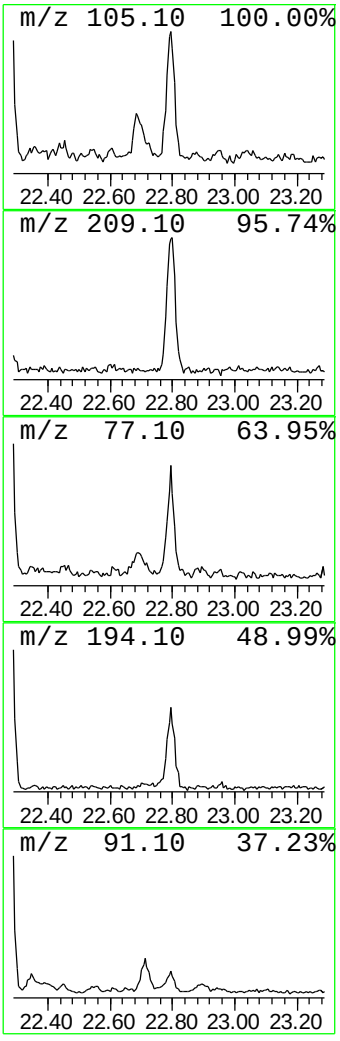
Peak Number 24 1-Heptanone, 1,3-diphenyl- Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.79	3.72 ng/ul	187080	Chrysene-d12	22.05

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Heptanone, 1,3-diphenyl-	266	C19H22O	001454-57-5	53
2	Carbazole, 2,4,7-trimethyl-	209	C15H15N	078787-89-0	53
3	9-Methoxyacridine	209	C14H11NO	010228-90-7	43
4	Carbazole, 1,3,4-trimethyl-	209	C15H15N	078787-82-3	43
5	1,3,5-Triphenyl-1,5-pentanedione	328	C23H20O2	006263-84-9	43





Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG080217\
Data File : BG028136.D
Acq On : 3 Aug 2017 11:33
Operator : SJ/JU
Sample : I4405-07
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : JJ2X4

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG080217.M
Quant Title : SVOA CALIBRATION

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

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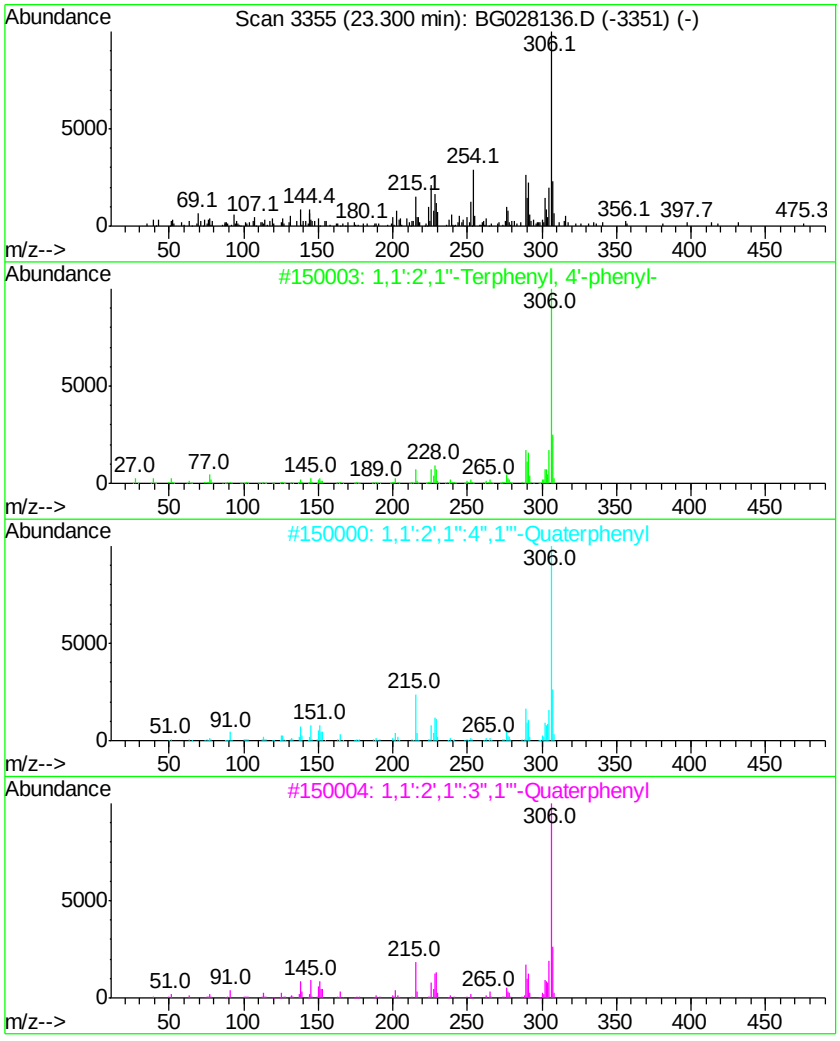
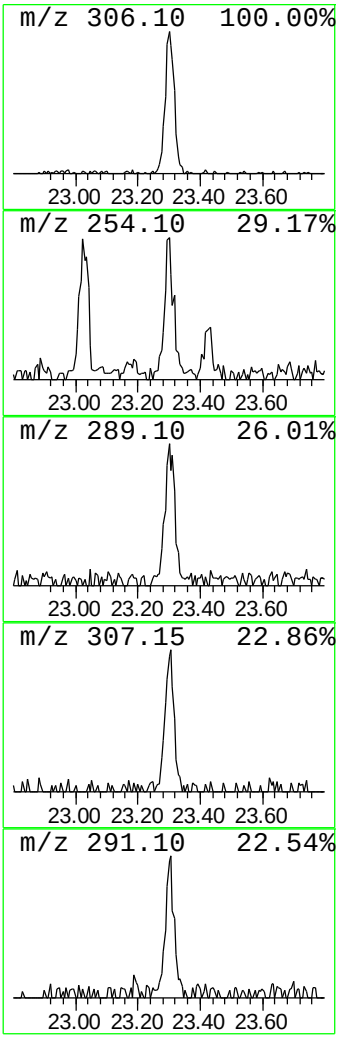
Peak Number 25 1,1':2',1''-Terphenyl, 4'-p... Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.30	3.24 ng/ul	163012	Chrysene-d12	22.05

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1':2',1''-Terphenyl, 4'-phenyl-	306	C24H18	001165-53-3	97
2	1,1':2',1'':4'',1'''-Quaterphenyl	306	C24H18	001165-58-8	93
3	1,1':2',1'':3'',1'''-Quaterphenyl	306	C24H18	001165-57-7	90
4	1,1':2',1'':2'',1'''-Quaterphenyl	306	C24H18	000641-96-3	86
5	1,1':3',1''-Terphenyl, 5'-phenyl-	306	C24H18	000612-71-5	86





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Data File : BG028136.D
Acq On : 3 Aug 2017 11:33
Operator : SJ/JU
Sample : I4405-07
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : JJ2X4

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG080217.M
Quant Title : SVOA CALIBRATION

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

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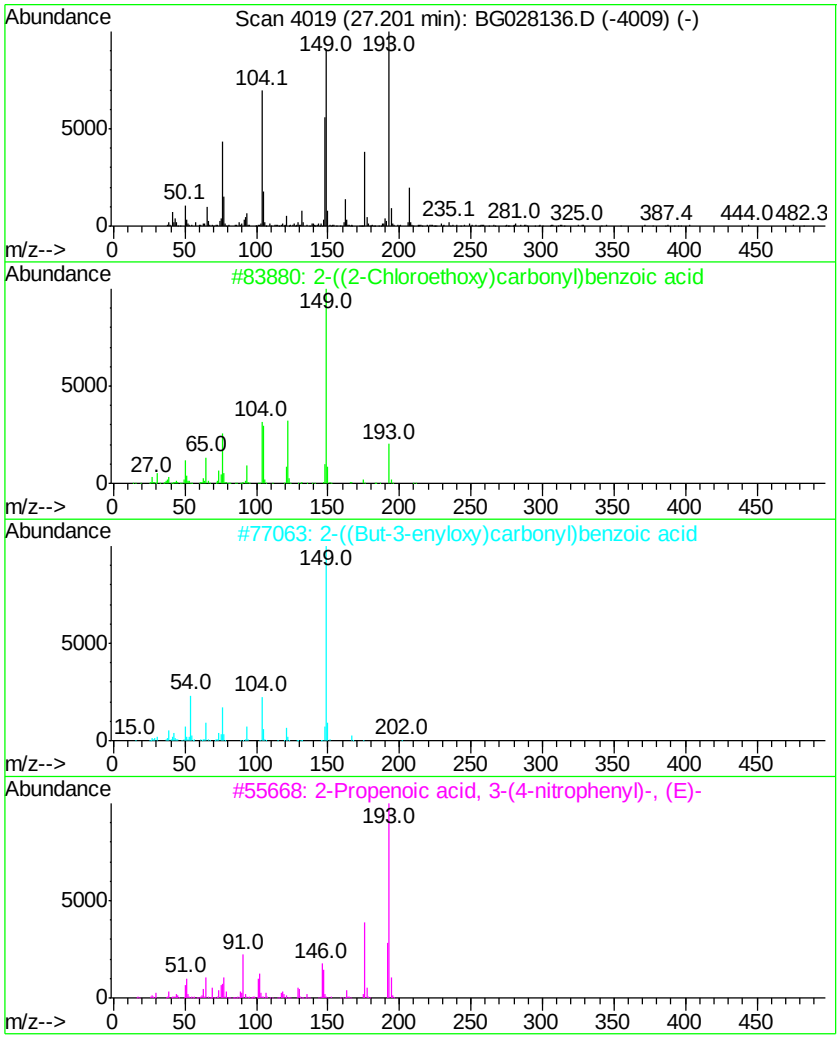
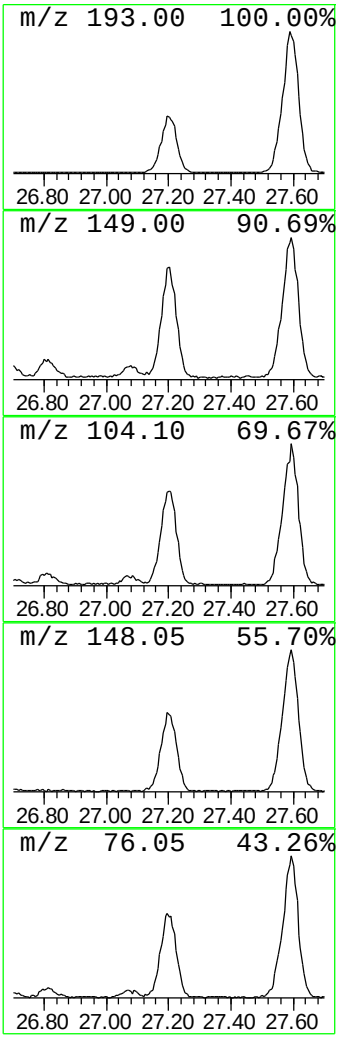
Peak Number 26 unknown-07 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
27.20	15.18 ng/ul	1130270	Perylene-d12	25.57

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-((2-Chloroethoxy)carbonyl)benz...	228	C10H9ClO4	1000373-53-5	38
2	2-((But-3-enyloxy)carbonyl)benzo...	220	C12H12O4	1000373-53-4	35
3	2-Propenoic acid, 3-(4-nitrophen...	193	C9H7NO4	000882-06-4	25
4	2-((Prop-2-ynyloxy)carbonyl)benz...	204	C11H8O4	1000373-53-6	25
5	Cyclohexanol, 4-phenyl-	176	C12H16O	005437-46-7	25





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Data File : BG028136.D
Acq On : 3 Aug 2017 11:33
Operator : SJ/JU
Sample : I4405-07
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : JJ2X4

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG080217.M
Quant Title : SVOA CALIBRATION

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

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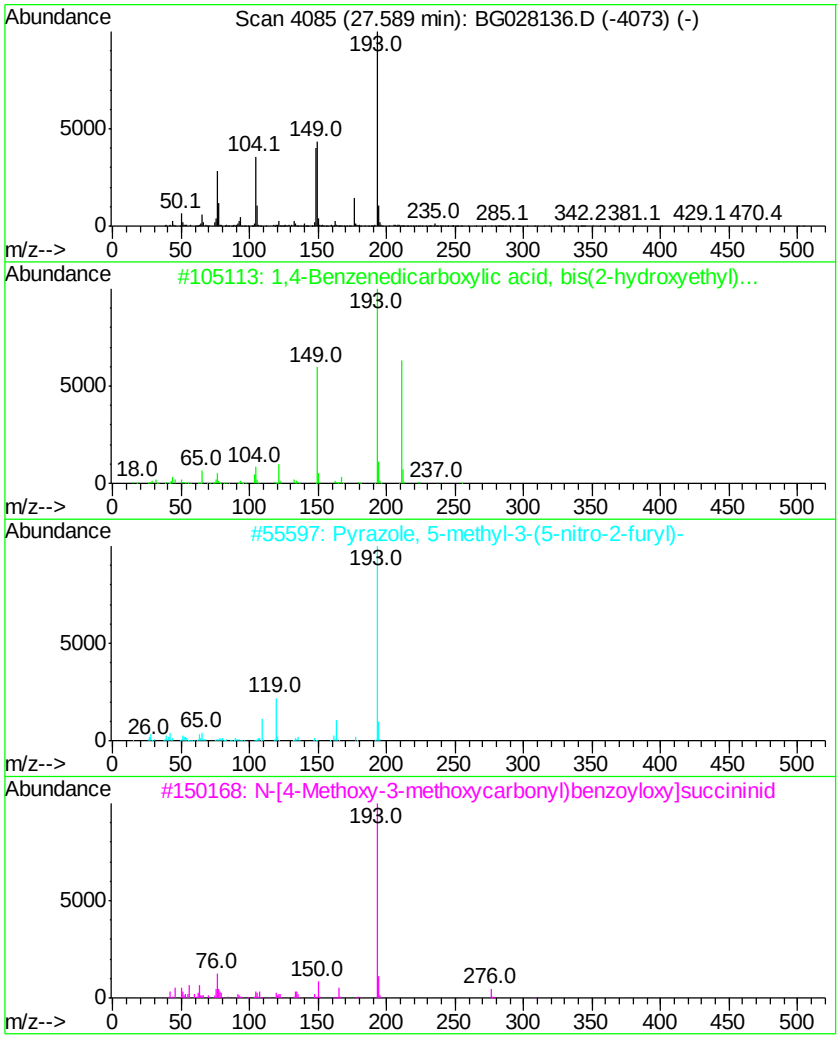
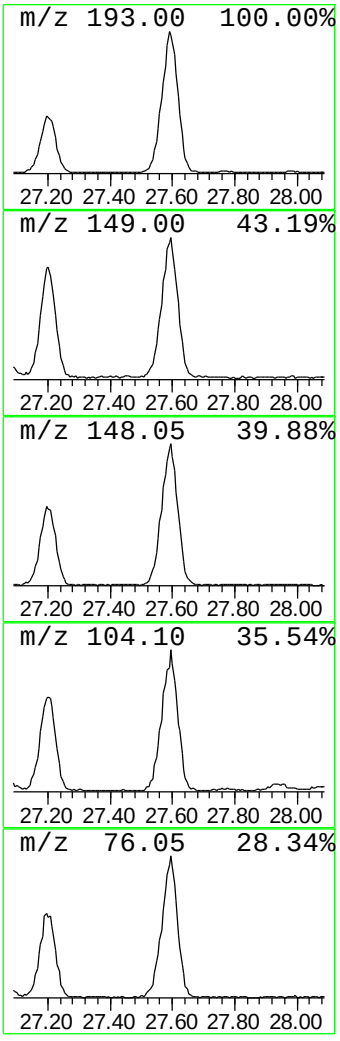
Peak Number 27 unknown-08 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
27.59	23.38 ng/ul	1740750	Perylene-d12	25.57

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,4-Benzenedicarboxylic acid, bi...	254	C12H14O6	000959-26-2	37
2	Pyrazole, 5-methyl-3-(5-nitro-2-...	193	C8H7N3O3	016239-90-0	25
3	N-[4-Methoxy-3-methoxycarbonyl)b...	307	C14H13NO7	541528-00-1	17
4	2-Aminobenzothiazol-6-carboxamide	193	C8H7N3OS	111962-90-4	16
5	Acridine, 9-methyl-	193	C14H11N	000611-64-3	12





Tentatively Identified Compound (LSC) summary

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG080217\
Data File : BG028136.D
Acq On : 3 Aug 2017 11:33
Operator : SJ/JU
Sample : I4405-07
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : JJ2X4

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG080217.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
Phthalic anhydride	12.93	9.6	ng/ul	219815	2	11.18	458295	20.0
Benzene, 1,1'-(1,...	16.50	11.9	ng/ul	507762	4	17.72	854817	20.0
1-Propene, 2-(3-m...	16.64	2.5	ng/ul	107202	4	17.72	854817	20.0
Benzene, 1,1'-(1-...	16.73	2.4	ng/ul	104018	4	17.72	854817	20.0
unknown-01	17.05	14.4	ng/ul	615290	4	17.72	854817	20.0
Naphthalene, 1,2,...	17.52	2.7	ng/ul	114504	4	17.72	854817	20.0
Benzene, 1,1'-(1,...	17.67	6.7	ng/ul	285047	4	17.72	854817	20.0
.beta.-Phenylprop...	18.11	6.8	ng/ul	292398	4	17.72	854817	20.0
2-(Hexyloxycarbon...	18.21	7.1	ng/ul	303745	4	17.72	854817	20.0
Phthalic acid, 3-...	18.33	2.0	ng/ul	86095	4	17.72	854817	20.0
unknown-02	18.41	5.1	ng/ul	218556	4	17.72	854817	20.0
n-Decanoic acid	18.58	5.9	ng/ul	251948	4	17.72	854817	20.0
1,2,4,8-Tetrameth...	19.08	4.3	ng/ul	182675	4	17.72	854817	20.0
Naphthalene, 2-(p...	19.86	9.6	ng/ul	412107	4	17.72	854817	20.0
unknown-03	21.05	2.2	ng/ul	108671	5	22.05	1006880	20.0
unknown-04	21.24	2.9	ng/ul	143696	5	22.05	1006880	20.0
unknown-05	21.36	3.8	ng/ul	190319	5	22.05	1006880	20.0
Pregna-5,17(20)-d...	21.61	8.1	ng/ul	408817	5	22.05	1006880	20.0
unknown-06	21.64	26.4	ng/ul	1331000	5	22.05	1006880	20.0
1-Propene, 3-(2-c...	22.28	33.6	ng/ul	1692720	5	22.05	1006880	20.0
1-Heptanone, 1,3-...	22.79	3.7	ng/ul	187080	5	22.05	1006880	20.0
1,1':2',1''-Terph...	23.30	3.2	ng/ul	163012	5	22.05	1006880	20.0
unknown-07	27.20	15.2	ng/ul	1130270	6	25.57	1488920	20.0
unknown-08	27.59	23.4	ng/ul	1740750	6	25.57	1488920	20.0