LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM062317\
Data File : BM010709.D
Acq On : 24 Jun 2017 02:38
Operator : SJ/JU
Sample : I3625-13ME 20X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F5C80ME

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 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:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM062017.M
Title : SVOA CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	7.463	771	778	784	rBV	300469	461756	17.43%	1.866%
2	7.828	833	840	846	rBB	30103	44467	1.68%	0.180%
3	7.986	860	867	873	rBB	50285	74345	2.81%	0.300%
4	10.222	1240	1247	1251	rBV	452777	766904	28.95%	3.098%
5	10.275	1251	1256	1264	rVB	1688771	2620874	98.95%	10.589%
6	10.386	1268	1275	1283	rVB	69720	124959	4.72%	0.505%
7	11.892	1525	1531	1538	rVB	519317	817568	30.87%	3.303%
8	12.016	1547	1552	1558	rVB2	19101	28744	1.09%	0.116%
9	12.121	1558	1570	1576	rBB	246542	390317	14.74%	1.577%
10	12.933	1703	1708	1715	rBB	123678	176844	6.68%	0.714%
11	13.133	1737	1742	1749	rBV3	36088	61177	2.31%	0.247%
12	13.263	1758	1764	1766	rBV2	63697	96866	3.66%	0.391%
13	13.427	1785	1792	1797	rBV2	104528	177205	6.69%	0.716%
14	13.480	1797	1801	1805	rVV	71417	96658	3.65%	0.391%
15	13.527	1805	1809	1814	rVB	37725	47109	1.78%	0.190%
16	13.663	1824	1832	1836	rBV4	40781	68286	2.58%	0.276%
17	13.798	1849	1855	1857	rBV2	34338	47703	1.80%	0.193%
18	13.827	1857	1860	1865	rVB3	40638	56668	2.14%	0.229%
19	14.110	1900	1908	1914	rBV2	760014	1203736	45.44%	4.863%
20	14.174	1914	1919	1927	rVV	521595	738676	27.89%	2.984%
21	14.239	1927	1930	1935	rVB	22758	30049	1.13%	0.121%
22	14.321	1940	1944	1954	rBV3	21940	35940	1.36%	0.145%
23	14.421	1956	1961	1967	rVB4	23107	34832	1.32%	0.141%
24	14.515	1969	1977	1984	rBV	429972	643544	24.30%	2.600%
25	14.592	1986	1990	1995	rBV2	22333	29478	1.11%	0.119%
26	14.733	2008	2014	2019	rBV4	18603	31950	1.21%	0.129%
27	14.792	2019	2024	2029	rVB4	20950	27103	1.02%	0.110%
28	15.110	2073	2078	2082	rBV2	59629	83098	3.14%	0.336%
29	15.168	2082	2088	2093	rVB	605260	819316	30.93%	3.310%
30	15.333	2112	2116	2120	rVB2	64007	88194	3.33%	0.356%
31	15.415	2127	2130	2134	rVB2	30535	36763	1.39%	0.149%
32	15.462	2134	2138	2146	rVB	86404	121217	4.58%	0.490%
33	15.609	2158	2163	2171	rVB	92970	176896	6.68%	0.715%
34	15.698	2171	2178	2185	rVB3	15637	30879	1.17%	0.125%

LSC Area Percent Report

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Operator : SJ/JU
Sample : I3625-13ME 20X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F5C80ME

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 1 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

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Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM062017.M
Title : SVOA CALIBRATION

35	15.968	2219	2224	2231	rBV4	23732	39874	1.51%	0.161%
36	16.174	2249	2259	2264	rBV3	60170	127986	4.83%	0.517%
37	16.221	2264	2267	2272	rVB3	22479	28753	1.09%	0.116%
38	16.321	2279	2284	2290	rVB3	24400	40819	1.54%	0.165%
39	16.409	2295	2299	2302	rVV3	23500	30559	1.15%	0.123%
40	16.445	2302	2305	2309	rVB4	20960	28853	1.09%	0.117%
41	16.604	2326	2332	2338	rBV3	32298	58796	2.22%	0.238%
42	16.680	2341	2345	2350	rVB	125151	170132	6.42%	0.687%
43	16.868	2371	2377	2380	rBV	1009349	1459611	55.11%	5.897%
44	16.909	2380	2384	2390	rVV	1980329	2648779	100.00%	10.701%
45	16.962	2390	2393	2395	rVV2	77335	96579	3.65%	0.390%
46	16.998	2395	2399	2405	rVB	265920	356310	13.45%	1.440%
47	17.056	2405	2409	2416	rBV4	38023	58046	2.19%	0.235%
48	17.274	2440	2446	2450	rBV	166302	221867	8.38%	0.896%
49	17.398	2462	2467	2470	rVB5	22763	30570	1.15%	0.124%
50	17.745	2516	2526	2529	rBV	122775	193398	7.30%	0.781%
51	17.792	2529	2534	2540	rVB	189533	275450	10.40%	1.113%
52	17.868	2540	2547	2553	rBV	73797	111660	4.22%	0.451%
53	17.939	2553	2559	2563	rVV2	217745	379830	14.34%	1.535%
54	17.974	2563	2565	2570	rVV2	69629	76073	2.87%	0.307%
55	18.050	2572	2578	2581	rVV3	17405	27022	1.02%	0.109%
56	18.256	2608	2613	2617	rVB	89809	117839	4.45%	0.476%
57	18.503	2649	2655	2659	rBV6	20426	29349	1.11%	0.119%
58	18.668	2679	2683	2689	rBV4	39576	74381	2.81%	0.301%
59	18.739	2690	2695	2698	rBV3	25681	36463	1.38%	0.147%
60	18.809	2704	2707	2716	rVB9	19571	46787	1.77%	0.189%
61	18.939	2723	2729	2735	rBV	1227798	1551433	58.57%	6.268%
62	19.039	2743	2746	2750	rVB5	20809	26786	1.01%	0.108%
63	19.180	2763	2770	2776	rVB3	35475	59303	2.24%	0.240%
64	19.280	2781	2787	2788	rBV2	271529	397714	15.01%	1.607%
65	19.303	2788	2791	2795	rVB	682646	821342	31.01%	3.318%
66	19.380	2800	2804	2810	rVB3	42915	72592	2.74%	0.293%
67	19.486	2816	2822	2829	rBV	52407	76453	2.89%	0.309%
68	19.550	2829	2833	2839	rVB2	36235	49098	1.85%	0.198%
69	19.633	2839	2847	2848	rBV3	43824	73660	2.78%	0.298%
70	19.692	2853	2857	2860	rVB	29100	34605	1.31%	0.140%
71	19.809	2873	2877	2882	rVB	152786	208289	7.86%	0.842%

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Instrument : BNA_M
ClientSampleId : F5C80ME

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 1 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
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Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM062017.M
Title : SVOA CALIBRATION

72	19.909	2889	2894	2898	rBV	192280	233560	8.82%	0.944%
73	19.962	2898	2903	2907	rVV3	61985	104599	3.95%	0.423%
74	20.150	2932	2935	2940	rBV5	20727	31618	1.19%	0.128%
75	20.339	2963	2967	2975	rBV4	27932	50489	1.91%	0.204%
76	20.521	2994	2998	3003	rBV5	20345	40246	1.52%	0.163%
77	20.727	3029	3033	3036	rBV	50733	69043	2.61%	0.279%
78	20.768	3036	3040	3043	rVV4	44245	66602	2.51%	0.269%
79	20.809	3043	3047	3055	rVB3	34797	75570	2.85%	0.305%
80	21.080	3086	3093	3097	rBV2	1377498	2002053	75.58%	8.089%
81	21.115	3097	3099	3106	rVB	202707	245058	9.25%	0.990%
82	21.233	3116	3119	3123	rVB4	39942	43060	1.63%	0.174%
83	21.391	3144	3146	3152	rVB	27699	33272	1.26%	0.134%
84	22.585	3346	3349	3352	rBV	43746	73806	2.79%	0.298%
85	23.033	3422	3425	3429	rBV5	25088	37677	1.42%	0.152%
86	23.133	3438	3442	3448	rVB2	50468	97164	3.67%	0.393%
87	23.221	3450	3457	3465	rBV	636048	1120535	42.30%	4.527%

Sum of corrected areas: 24751534

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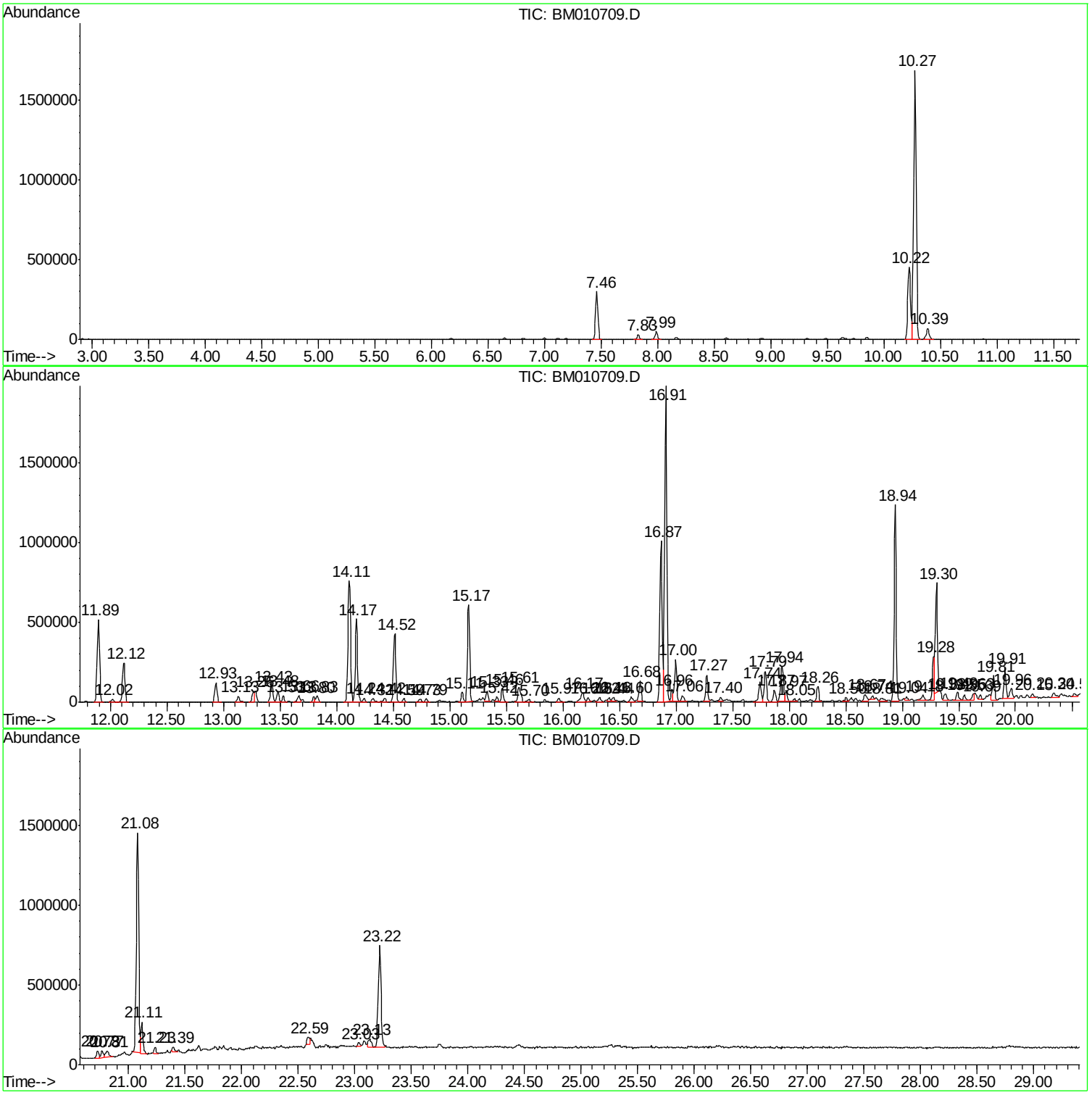
Instrument :
BNA_M
ClientSampleId :
F5C80ME

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM062017.M
Quant Title : SVOA CALIBRATION

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



Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM062317\
Data File : BM010709.D
Acq On : 24 Jun 2017 02:38
Operator : SJ/JU
Sample : I3625-13ME 20X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F5C80ME

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM062017.M
Quant Title : SVOA CALIBRATION

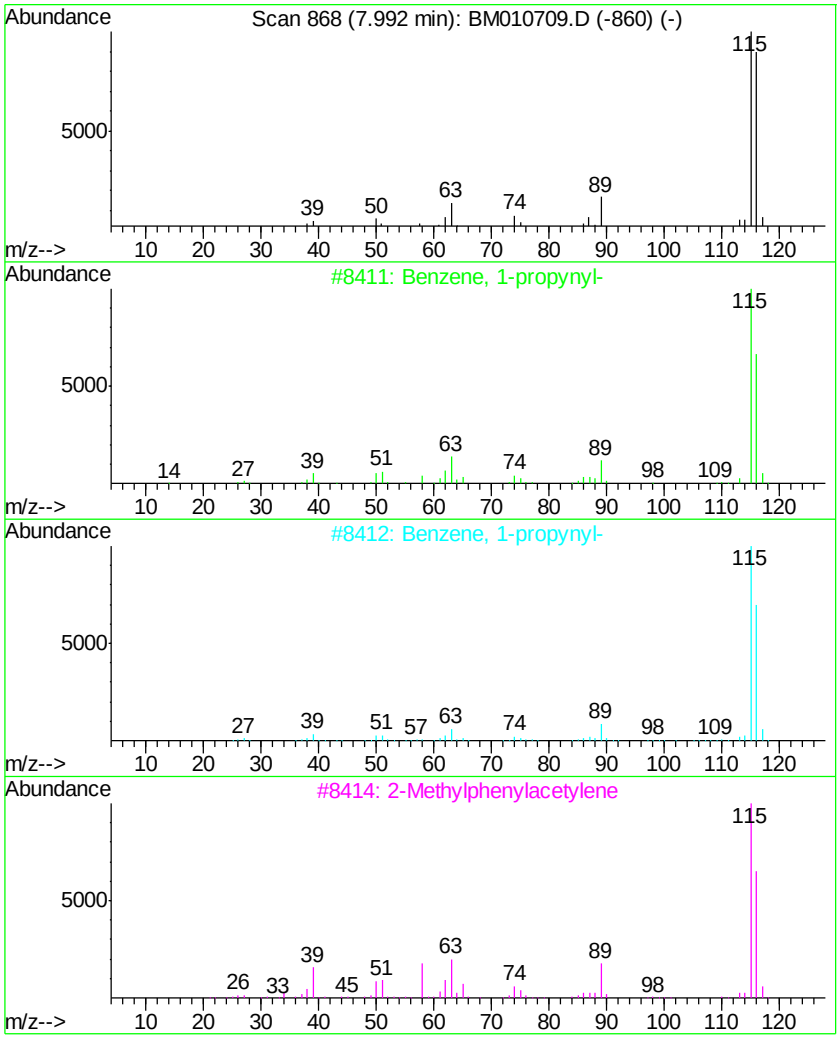
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

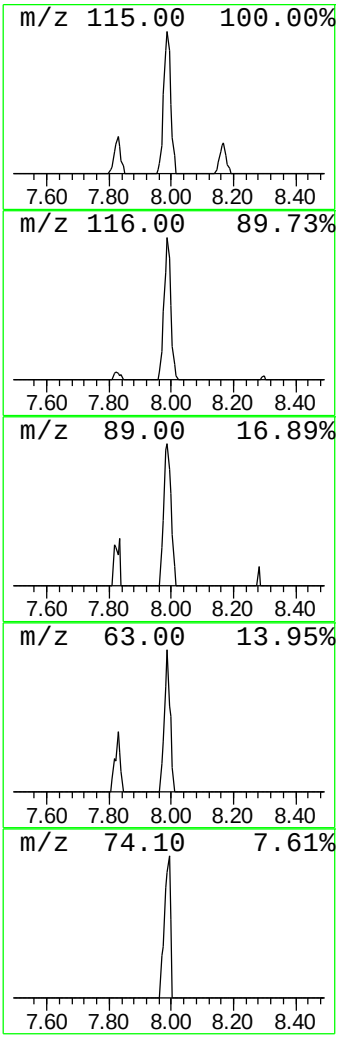
****

Peak Number 1 Benzene, 1-propynyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.99	3.22 ng/ul	74345	1,4-Dichlorobenzene-d4	7.46

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-propynyl-	116	C9H8	000673-32-5	91
2	Benzene, 1-propynyl-	116	C9H8	000673-32-5	90
3	2-Methylphenylacetylene	116	C9H8	000766-47-2	90
4	1-Propyne, 3-phenyl-	116	C9H8	010147-11-2	90
5	Benzene, 1-propynyl-	116	C9H8	000673-32-5	90





Library Search Compound Report

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Operator : SJ/JU
Sample : I3625-13ME 20X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F5C80ME

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM062017.M
Quant Title : SVOA CALIBRATION

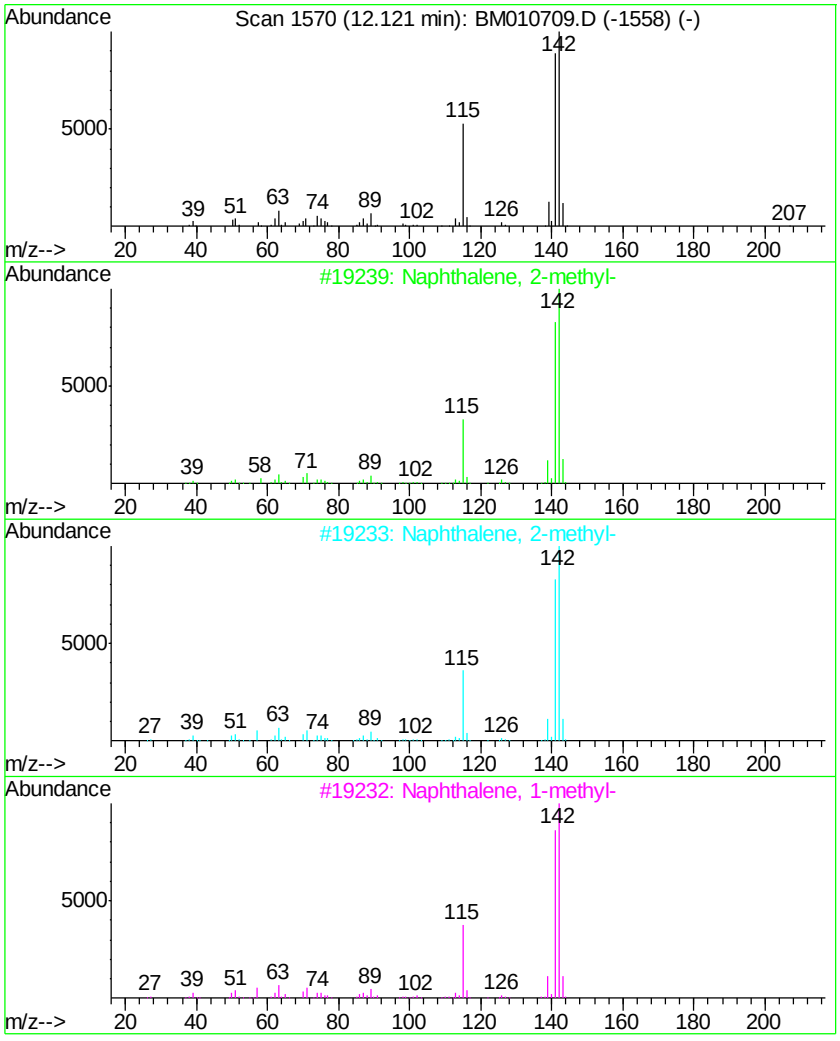
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

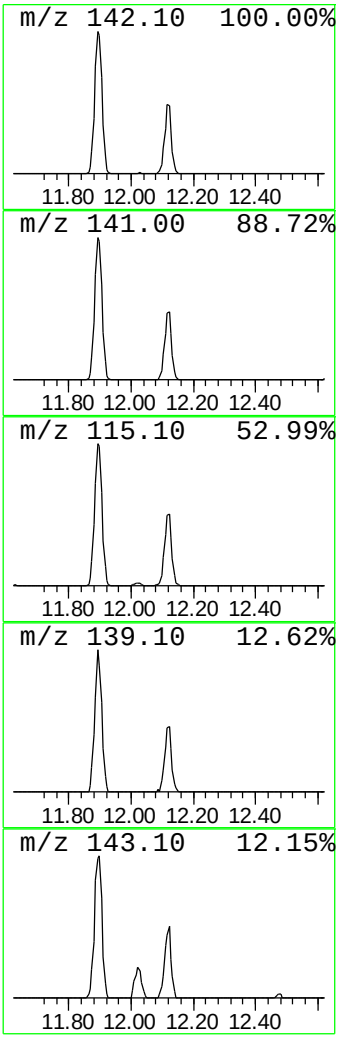
****

Peak Number 2 Naphthalene, 1-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.12	10.18 ng/ul	390317	Naphthalene-d8	10.22

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
2	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
3	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
4	Benzocycloheptatriene	142	C11H10	000264-09-5	94
5	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	94





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Instrument : BNA_M
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Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM062017.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
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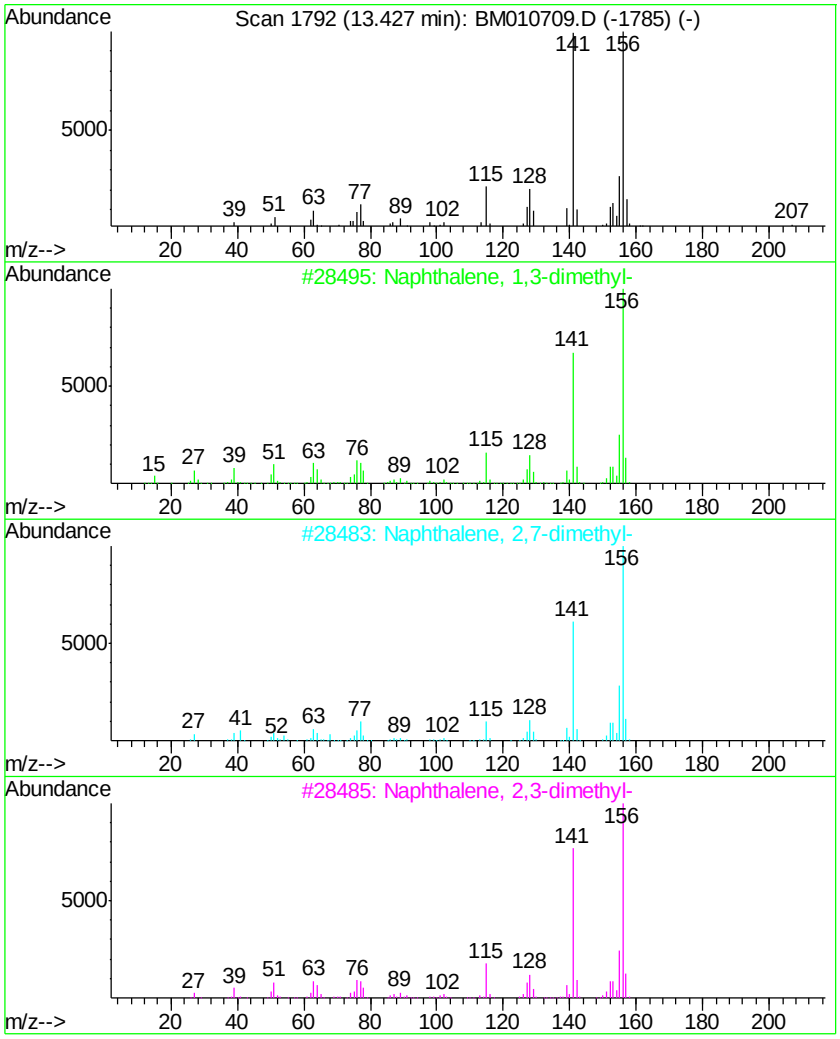
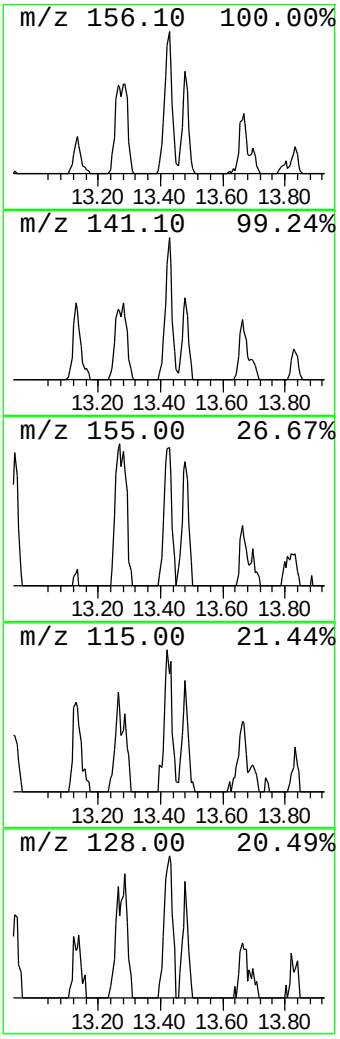
****

Peak Number 3 Naphthalene, 1,3-dimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.43	2.94 ng/ul	177205	Acenaphthene-d10	14.11

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	97
2	Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	97
3	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
4	Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	97
5	Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	96





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Quant Title : SVOA CALIBRATION

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

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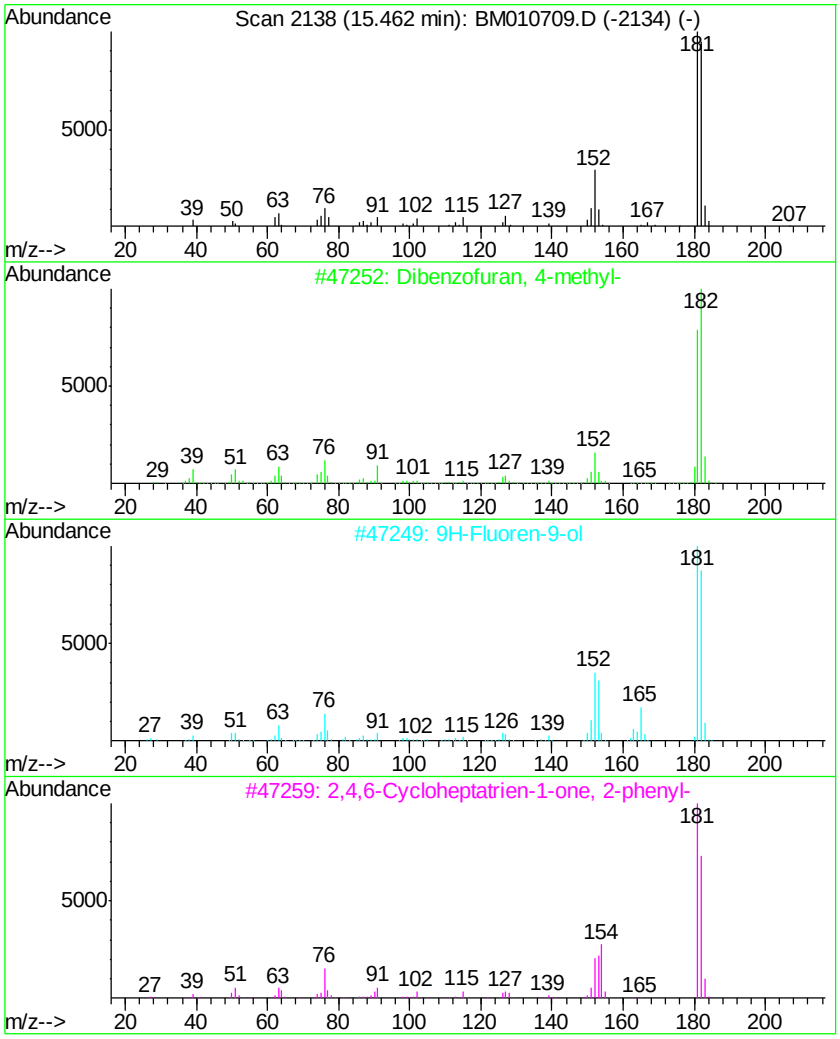
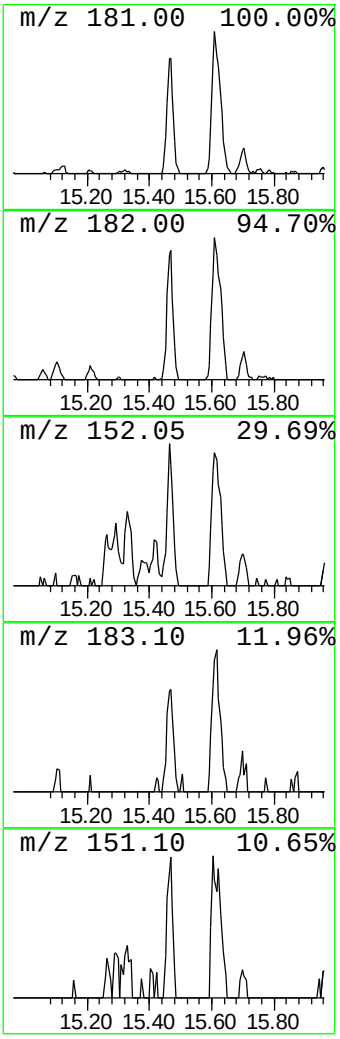
Peak Number 4 Dibenzofuran, 4-methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.46	2.01 ng/ul	121217	Acenaphthene-d10	14.11

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dibenzofuran, 4-methyl-	182	C13H10O	007320-53-8	91
2	9H-Fluoren-9-ol	182	C13H10O	001689-64-1	87
3	2,4,6-Cycloheptatrien-1-one, 2-p...	182	C13H10O	014562-09-5	87
4	9H-Fluoren-9-ol	182	C13H10O	001689-64-1	86
5	[1,1'-Biphenyl]-4-carboxaldehyde	182	C13H10O	003218-36-8	68





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Sample : I3625-13ME 20X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F5C80ME

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM062017.M
Quant Title : SVOA CALIBRATION

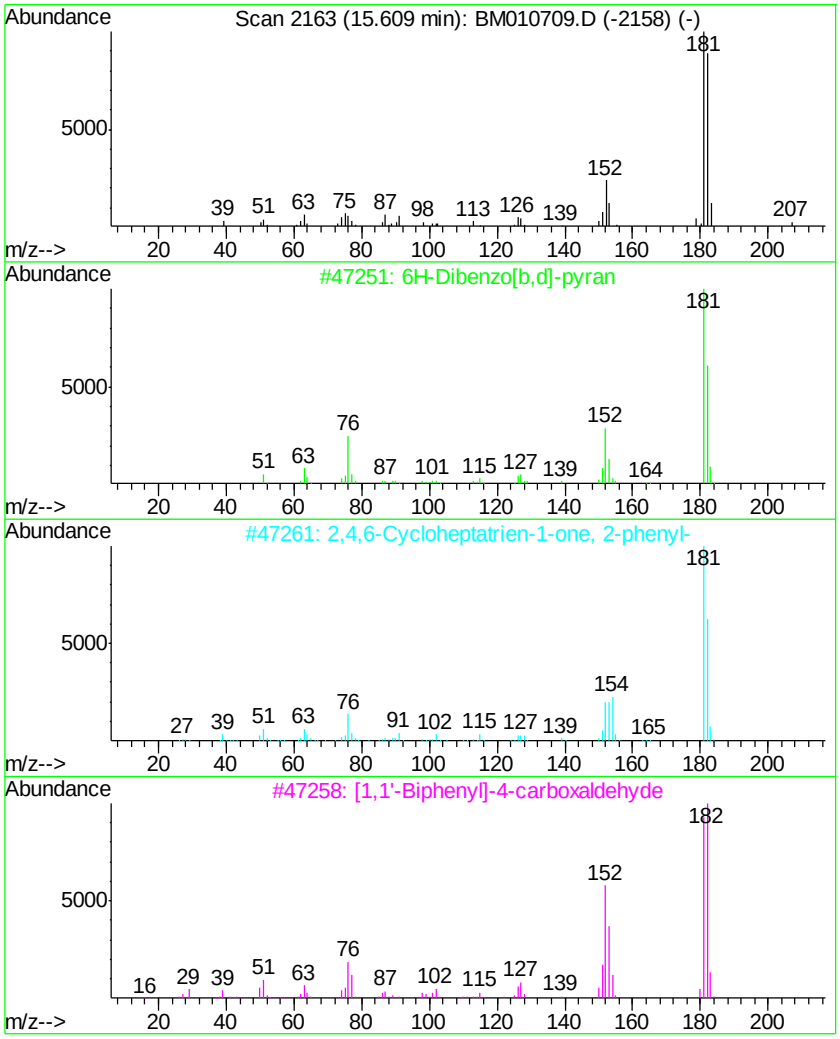
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

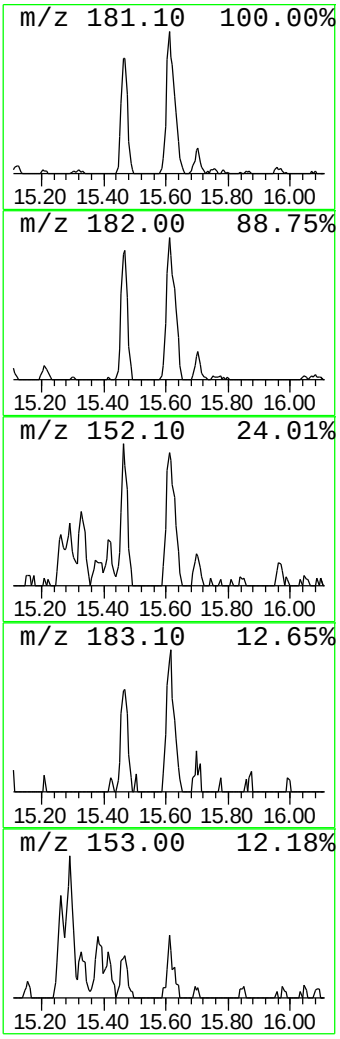
****

Peak Number 5 6H-Dibenzo[b,d]-pyran Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.61	2.42 ng/ul	176896	Phenanthrene-d10	16.87

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	6H-Dibenzo[b,d]-pyran	182	C13H10O	000229-95-8	91
2	2,4,6-Cycloheptatrien-1-one, 2-p...	182	C13H10O	014562-09-5	90
3	[1,1'-Biphenyl]-4-carboxaldehyde	182	C13H10O	003218-36-8	86
4	9H-Fluoren-9-ol	182	C13H10O	001689-64-1	86
5	2,4,6-Cycloheptatrien-1-one, 2-p...	182	C13H10O	014562-09-5	86





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Quant Title : SVOA CALIBRATION

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

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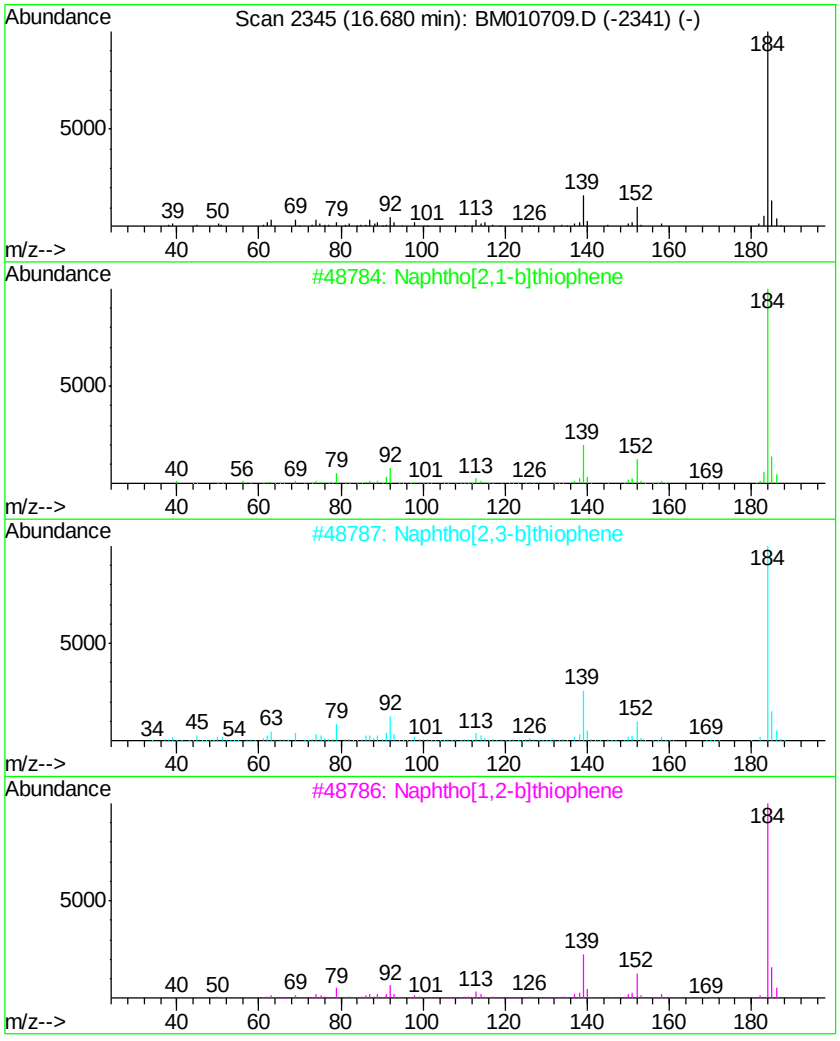
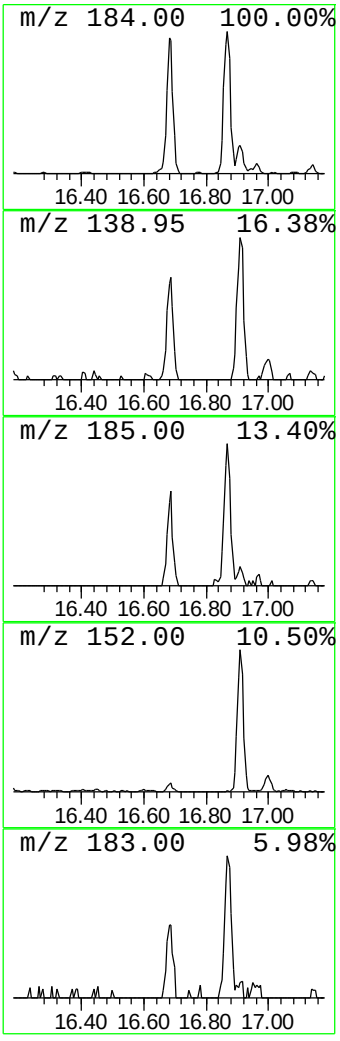
Peak Number 6 Naphtho[2,1-b]thiophene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.68	2.33 ng/ul	170132	Phenanthrene-d10	16.87

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphtho[2,1-b]thiophene	184	C12H8S	000233-02-3	97
2	Naphtho[2,3-b]thiophene	184	C12H8S	000268-77-9	95
3	Naphtho[1,2-b]thiophene	184	C12H8S	000234-41-3	94
4	Dibenzothiophene	184	C12H8S	000132-65-0	93
5	Dibenzothiophene	184	C12H8S	000132-65-0	93





Data Path : Z:\HPCHEM1\BNA_M\DATA\BM062317\
Data File : BM010709.D
Acq On : 24 Jun 2017 02:38
Operator : SJ/JU
Sample : I3625-13ME 20X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F5C80ME

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM062017.M
Quant Title : SVOA CALIBRATION

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

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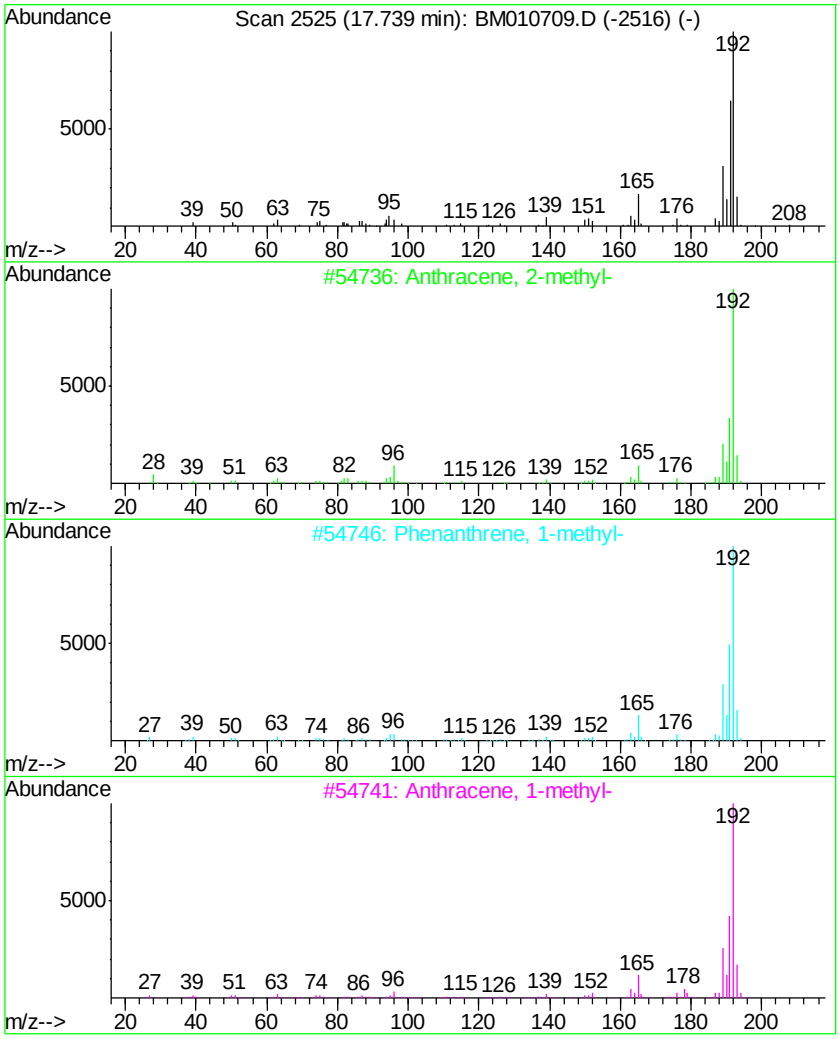
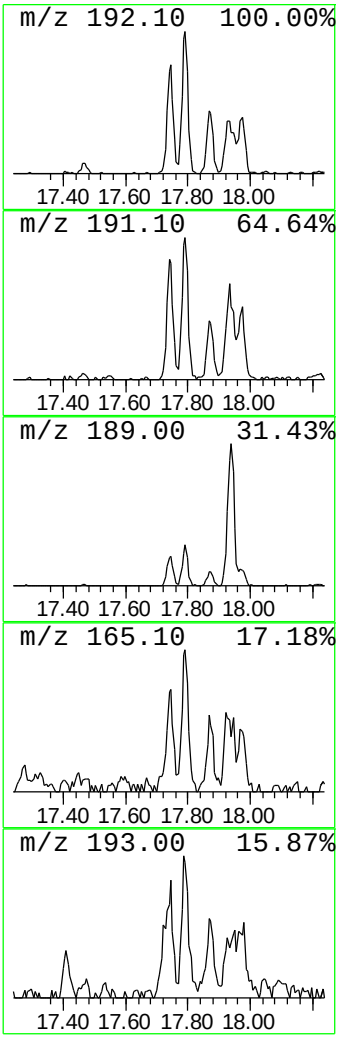
Peak Number 7 Anthracene, 2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.74	2.65 ng/ul	193398	Phenanthrene-d10	16.87

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Anthracene, 2-methyl-	192	C15H12	000613-12-7	93
2	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	93
3	Anthracene, 1-methyl-	192	C15H12	000610-48-0	91
4	Anthracene, 1-methyl-	192	C15H12	000610-48-0	91
5	Anthracene, 2-methyl-	192	C15H12	000613-12-7	91





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Data File : BM010709.D
Acq On : 24 Jun 2017 02:38
Operator : SJ/JU
Sample : I3625-13ME 20X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F5C80ME

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM062017.M
Quant Title : SVOA CALIBRATION

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

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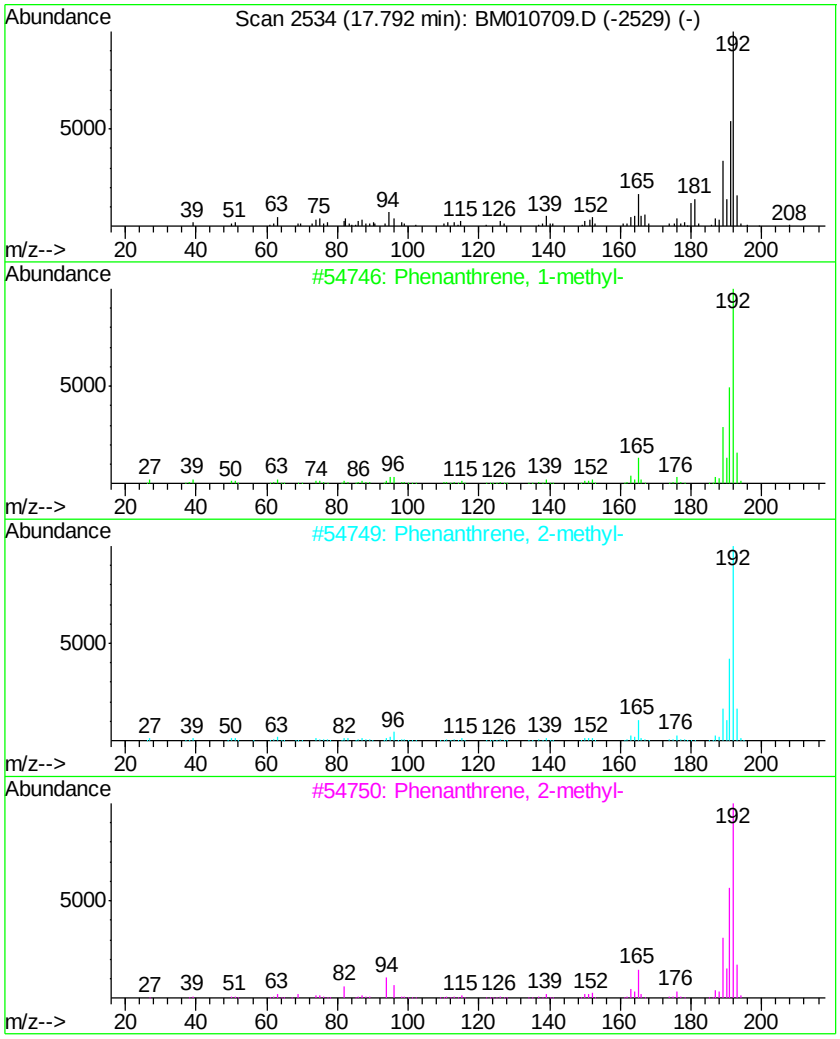
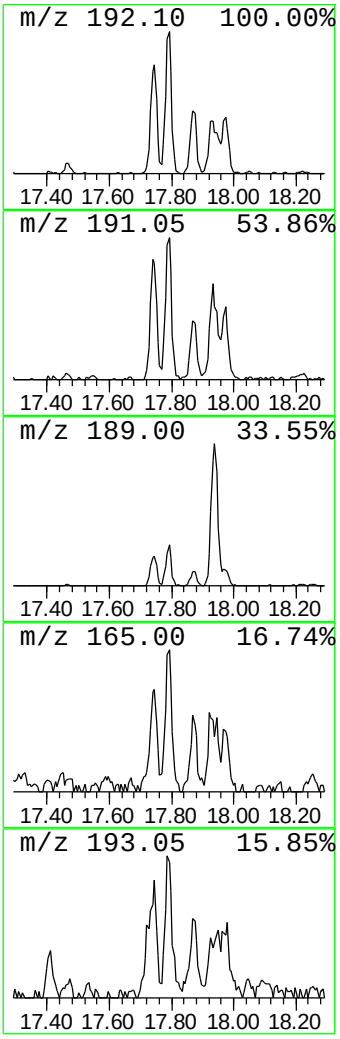
Peak Number 8 Phenanthrene, 1-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.79	3.77 ng/ul	275450	Phenanthrene-d10	16.87

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	95
2	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	95
3	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	95
4	Anthracene, 9-methyl-	192	C15H12	000779-02-2	95
5	Anthracene, 2-methyl-	192	C15H12	000613-12-7	95





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Data File : BM010709.D
Acq On : 24 Jun 2017 02:38
Operator : SJ/JU
Sample : I3625-13ME 20X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F5C80ME

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM062017.M
Quant Title : SVOA CALIBRATION

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

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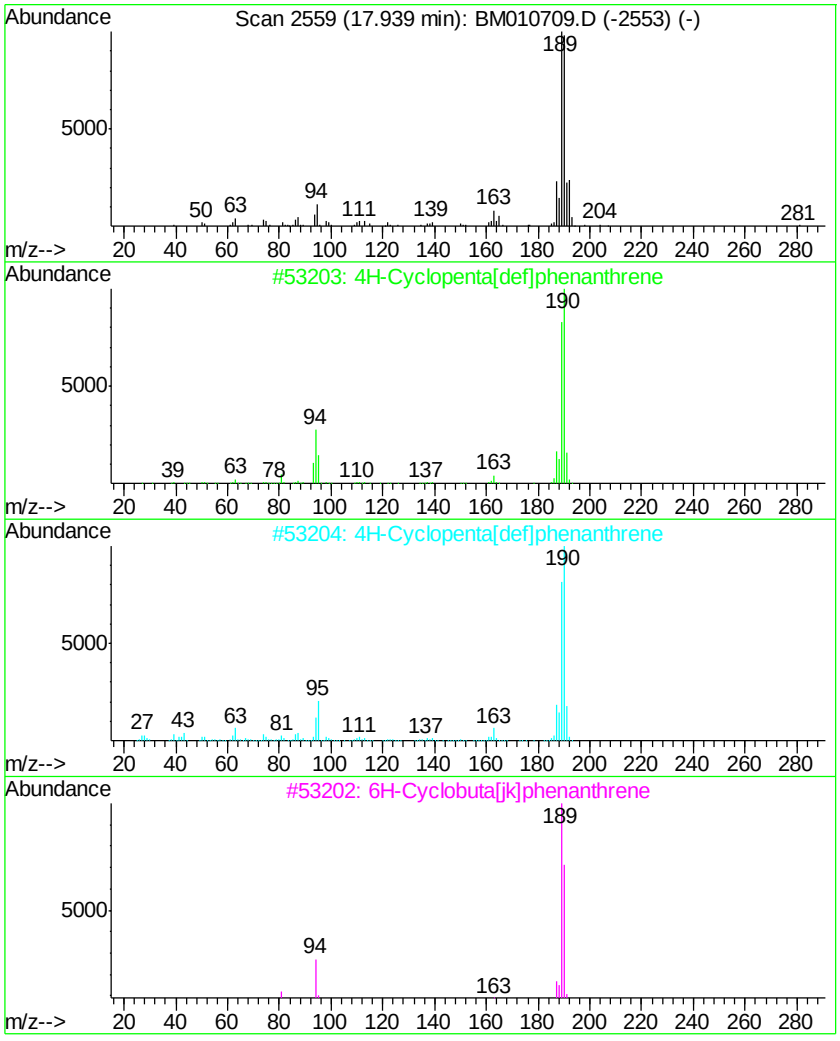
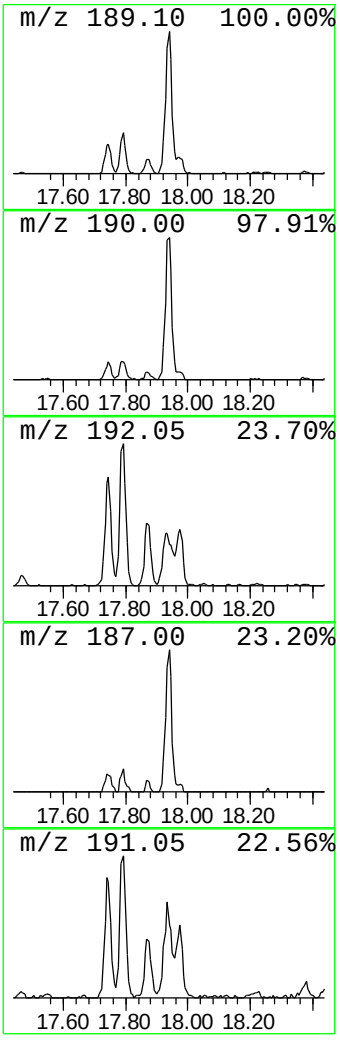
Peak Number 9 4H-Cyclopenta[def]phenanthrene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.94	5.20 ng/ul	379830	Phenanthrene-d10	16.87

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	87
2	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	81
3	6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	64
4	Methyl diselenide	190	C2H6Se2	007101-31-7	55
5	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	50





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Operator : SJ/JU
Sample : I3625-13ME 20X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F5C80ME

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM062017.M
Quant Title : SVOA CALIBRATION

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

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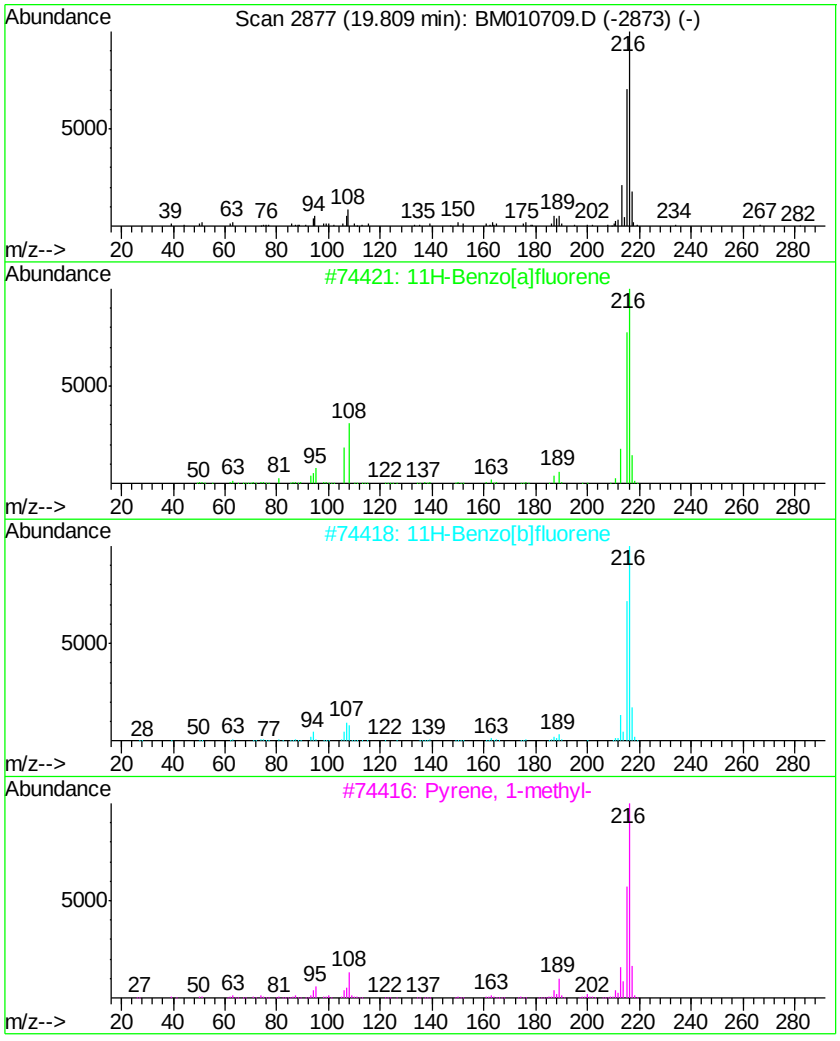
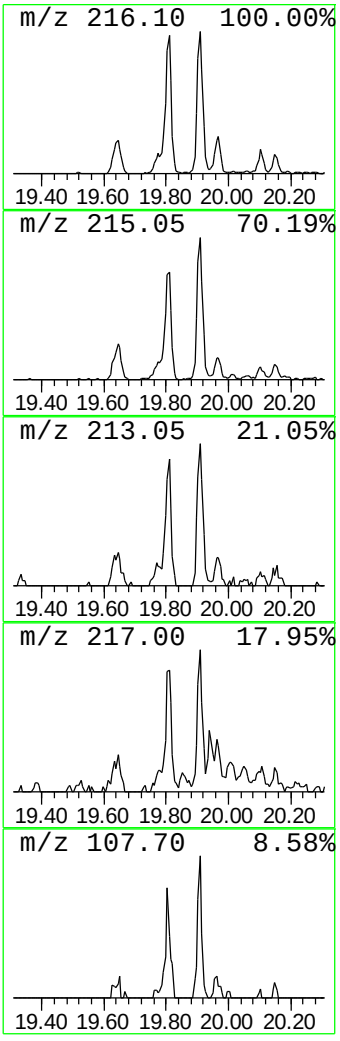
Peak Number 10 11H-Benzo[a]fluorene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.81	2.08 ng/ul	208289	Chrysene-d12	21.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	90
2	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	90
3	Pyrene, 1-methyl-	216	C17H12	002381-21-7	90
4	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	90
5	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	87





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Operator : SJ/JU
Sample : I3625-13ME 20X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F5C80ME

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM062017.M
Quant Title : SVOA CALIBRATION

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

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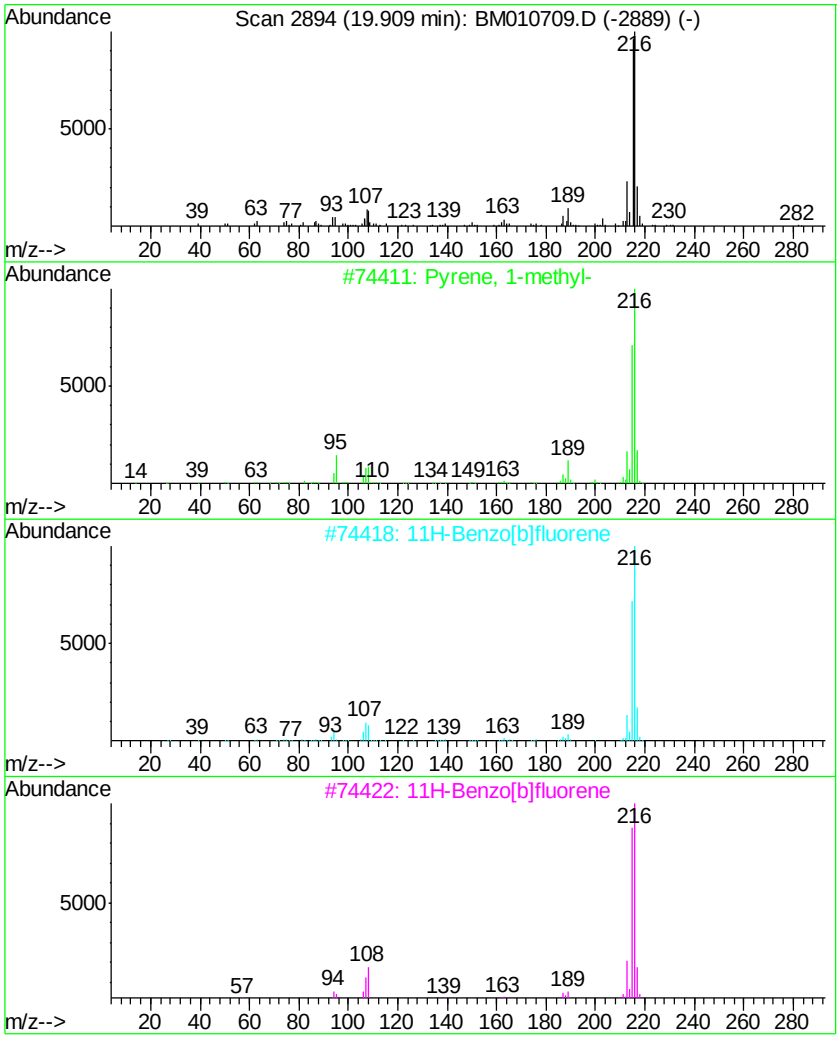
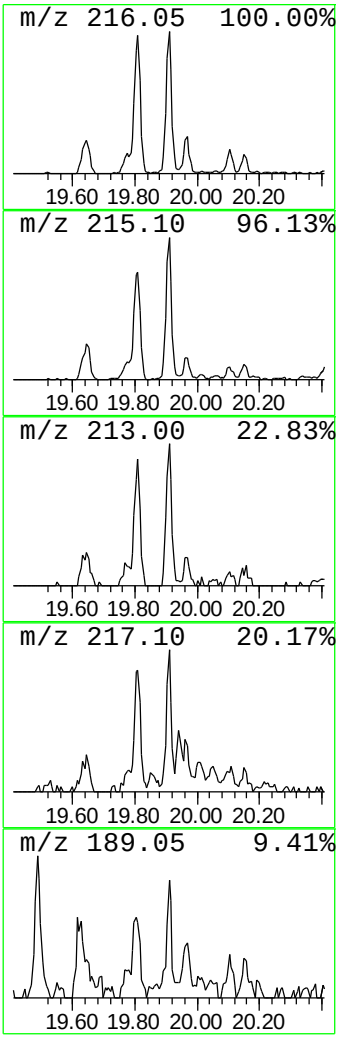
Peak Number 11 Pyrene, 1-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.91	2.33 ng/ul	233560	Chrysene-d12	21.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pyrene, 1-methyl-	216	C17H12	002381-21-7	94
2	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	90
3	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	87
4	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	83
5	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	80





Tentatively Identified Compound (LSC) summary

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM062317\
Data File : BM010709.D
Acq On : 24 Jun 2017 02:38
Operator : SJ/JU
Sample : I3625-13ME 20X
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : F5C80ME

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM062017.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
Benzene, 1-propynyl-	7.99	3.2	ng/ul	74345	1	7.46	461756	20.0
Naphthalene, 1-me...	12.12	10.2	ng/ul	390317	2	10.22	766904	20.0
Naphthalene, 1,3-...	13.43	2.9	ng/ul	177205	3	14.11	1203740	20.0
Dibenzofuran, 4-m...	15.46	2.0	ng/ul	121217	3	14.11	1203740	20.0
6H-Dibenzo[b,d]-p...	15.61	2.4	ng/ul	176896	4	16.87	1459610	20.0
Naphtho[2,1-b]thi...	16.68	2.3	ng/ul	170132	4	16.87	1459610	20.0
Anthracene, 2-met...	17.74	2.6	ng/ul	193398	4	16.87	1459610	20.0
Phenanthrene, 1-m...	17.79	3.8	ng/ul	275450	4	16.87	1459610	20.0
4H-Cyclopenta[def...	17.94	5.2	ng/ul	379830	4	16.87	1459610	20.0
11H-Benzo[a]fluorene	19.81	2.1	ng/ul	208289	5	21.08	2002050	20.0
Pyrene, 1-methyl-	19.91	2.3	ng/ul	233560	5	21.08	2002050	20.0