LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM100917\
Data File : BM012035.D
Acq On : 10 Oct 2017 04:40
Operator : SJ/JU
Sample : I5533-07
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : DUP-1

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-BM100317.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	7.004	659	666	680	rBV	861783	1645079	11.95%	0.958%
2	7.175	688	695	705	rBV	674502	1112525	8.08%	0.648%
3	7.375	722	729	739	rBV	943483	1615630	11.74%	0.941%
4	7.846	798	809	822	rBV	1107196	1887611	13.71%	1.099%
5	8.551	922	929	948	rBV	971688	1766006	12.83%	1.028%
6	9.010	1000	1007	1021	rBV	918672	1604189	11.65%	0.934%
7	9.734	1122	1130	1140	rBV	776045	1382794	10.04%	0.805%
8	10.269	1214	1221	1234	rBV	1216197	2200154	15.98%	1.281%
9	10.645	1274	1285	1291	rBV	1594509	2765286	20.09%	1.610%
10	10.786	1299	1309	1326	rBV	780765	1601828	11.64%	0.933%
11	13.898	1828	1838	1842	rVV	2420039	3463041	25.16%	2.016%
12	13.945	1842	1846	1852	rVB	576822	822996	5.98%	0.479%
13	14.186	1880	1887	1900	rBV	2769881	4296573	31.21%	2.501%
14	14.492	1927	1939	1945	rBV2	2508889	3919046	28.47%	2.281%
15	14.557	1945	1950	1957	rVB	234453	391651	2.84%	0.228%
16	14.692	1968	1973	1985	rBV	519005	1089692	7.92%	0.634%
17	14.798	1986	1991	1998	rVV5	58982	148248	1.08%	0.086%
18	14.892	1999	2007	2016	rVB	156791	333108	2.42%	0.194%
19	15.321	2076	2080	2085	rVB2	92631	139098	1.01%	0.081%
20	15.486	2099	2108	2113	rBV	3519898	5265017	38.24%	3.065%
21	15.539	2113	2117	2124	rVB	240657	378729	2.75%	0.220%
22	15.610	2124	2129	2135	rBV	831593	1221554	8.87%	0.711%
23	15.710	2141	2146	2156	rVB6	58111	170084	1.24%	0.099%
24	15.833	2162	2167	2173	rVB	117707	190213	1.38%	0.111%
25	15.980	2188	2192	2200	rVB	110460	210011	1.53%	0.122%
26	16.186	2222	2227	2229	rBV	124158	172814	1.26%	0.101%
27	16.521	2276	2284	2285	rBV4	87826	153420	1.11%	0.089%
28	16.539	2285	2287	2292	rVB3	100295	151547	1.10%	0.088%
29	16.768	2318	2326	2329	rBV3	138221	280861	2.04%	0.164%
30	16.892	2343	2347	2354	rVB2	134692	217192	1.58%	0.126%
31	16.974	2355	2361	2367	rVV4	182618	395909	2.88%	0.230%
32	17.057	2371	2375	2382	rVB	207472	338425	2.46%	0.197%
33	17.239	2399	2406	2409	rBV	3349890	4971246	36.11%	2.894%
34	17.280	2409	2413	2418	rVV	4543278	6153927	44.70%	3.583%

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Operator : SJ/JU
Sample : I5533-07
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : DUP-1

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-BM100317.M
Title : SVOA CALIBRATION

35	17.333	2418	2422	2426	rVV2	3851760	5754074	41.80%	3.350%
36	17.368	2426	2428	2434	rVB	957783	1165132	8.46%	0.678%
37	17.639	2468	2474	2483	rBV3	167492	384502	2.79%	0.224%
38	17.751	2490	2493	2500	rBV4	127413	205739	1.49%	0.120%
39	17.815	2500	2504	2513	rVB3	94908	170100	1.24%	0.099%
40	18.115	2549	2555	2559	rBV2	2177806	3264320	23.71%	1.900%
41	18.157	2559	2562	2570	rVV	1001254	1448781	10.52%	0.843%
42	18.239	2571	2576	2581	rVV	337192	483127	3.51%	0.281%
43	18.310	2581	2588	2591	rVV3	868669	1673132	12.15%	0.974%
44	18.339	2591	2593	2599	rVB	517775	670456	4.87%	0.390%
45	18.410	2601	2605	2610	rBV4	172926	270705	1.97%	0.158%
46	18.610	2634	2639	2643	rBV	573700	769781	5.59%	0.448%
47	18.657	2643	2647	2653	rVV	354484	495774	3.60%	0.289%
48	18.857	2677	2681	2685	rVV	215506	279690	2.03%	0.163%
49	18.904	2686	2689	2692	rVV	181638	212967	1.55%	0.124%
50	18.945	2692	2696	2699	rVB	165451	189783	1.38%	0.110%
51	19.027	2705	2710	2716	rBV2	488220	1041087	7.56%	0.606%
52	19.092	2718	2721	2724	rVB3	119420	168378	1.22%	0.098%
53	19.168	2729	2734	2742	rBV3	357176	691914	5.03%	0.403%
54	19.292	2749	2755	2761	rVB	8331213	11869798	86.22%	6.910%
55	19.351	2762	2765	2768	rVB2	161045	186761	1.36%	0.109%
56	19.392	2768	2772	2778	rBV2	1954300	2423539	17.60%	1.411%
57	19.539	2794	2797	2806	rVB2	309129	558279	4.06%	0.325%
58	19.627	2806	2812	2815	rVV3	6448647	10353322	75.21%	6.027%
59	19.657	2815	2817	2824	rVV	7776614	9297854	67.54%	5.413%
60	19.727	2824	2829	2835	rVB3	454632	778970	5.66%	0.453%
61	19.833	2844	2847	2853	rVB	369734	505429	3.67%	0.294%
62	19.898	2853	2858	2865	rVB2	250437	426592	3.10%	0.248%
63	19.974	2866	2871	2879	rBV3	524151	1366349	9.93%	0.795%
64	20.115	2889	2895	2896	rBV2	465271	734501	5.34%	0.428%
65	20.145	2897	2900	2905	rVB	995224	1415558	10.28%	0.824%
66	20.251	2914	2918	2921	rBV	645570	1096034	7.96%	0.638%
67	20.304	2923	2927	2931	rVV2	767945	1433359	10.41%	0.834%
68	20.445	2948	2951	2954	rBV	374519	392966	2.85%	0.229%
69	20.492	2956	2959	2962	rVV	422997	517030	3.76%	0.301%
70	20.533	2963	2966	2971	rVB	1134507	1339909	9.73%	0.780%
71	20.645	2982	2985	2990	rVB	290347	352377	2.56%	0.205%

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Misc :
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Instrument : BNA_M
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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-BM100317.M
Title : SVOA CALIBRATION

72	20.898	3024	3028	3033	rBV	666589	971501	7.06%	0.566%
73	21.098	3058	3062	3066	rVV3	1466946	1976855	14.36%	1.151%
74	21.139	3066	3069	3074	rVV2	610876	936278	6.80%	0.545%
75	21.192	3075	3078	3084	rVB	718416	938758	6.82%	0.546%
76	21.409	3109	3115	3119	rBV2	6800113	13766649	100.00%	8.014%
77	21.451	3119	3122	3131	rVB	4383635	5814104	42.23%	3.385%
78	21.568	3139	3142	3149	rVB	659628	852943	6.20%	0.497%
79	21.974	3208	3211	3215	rBV2	591500	780074	5.67%	0.454%
80	23.033	3385	3391	3396	rBV	2918379	6551939	47.59%	3.814%
81	23.203	3417	3420	3427	rVB	539494	895080	6.50%	0.521%
82	23.533	3470	3476	3479	rBV	1457623	2860484	20.78%	1.665%
83	23.580	3479	3484	3489	rVV2	3213193	6422403	46.65%	3.739%
84	23.633	3489	3493	3500	rVB	2458364	4438871	32.24%	2.584%
85	23.733	3504	3510	3515	rBV	2638225	4997209	36.30%	2.909%
86	26.097	3906	3912	3922	rVB	1016962	2729955	19.83%	1.589%

Sum of corrected areas: 171776676

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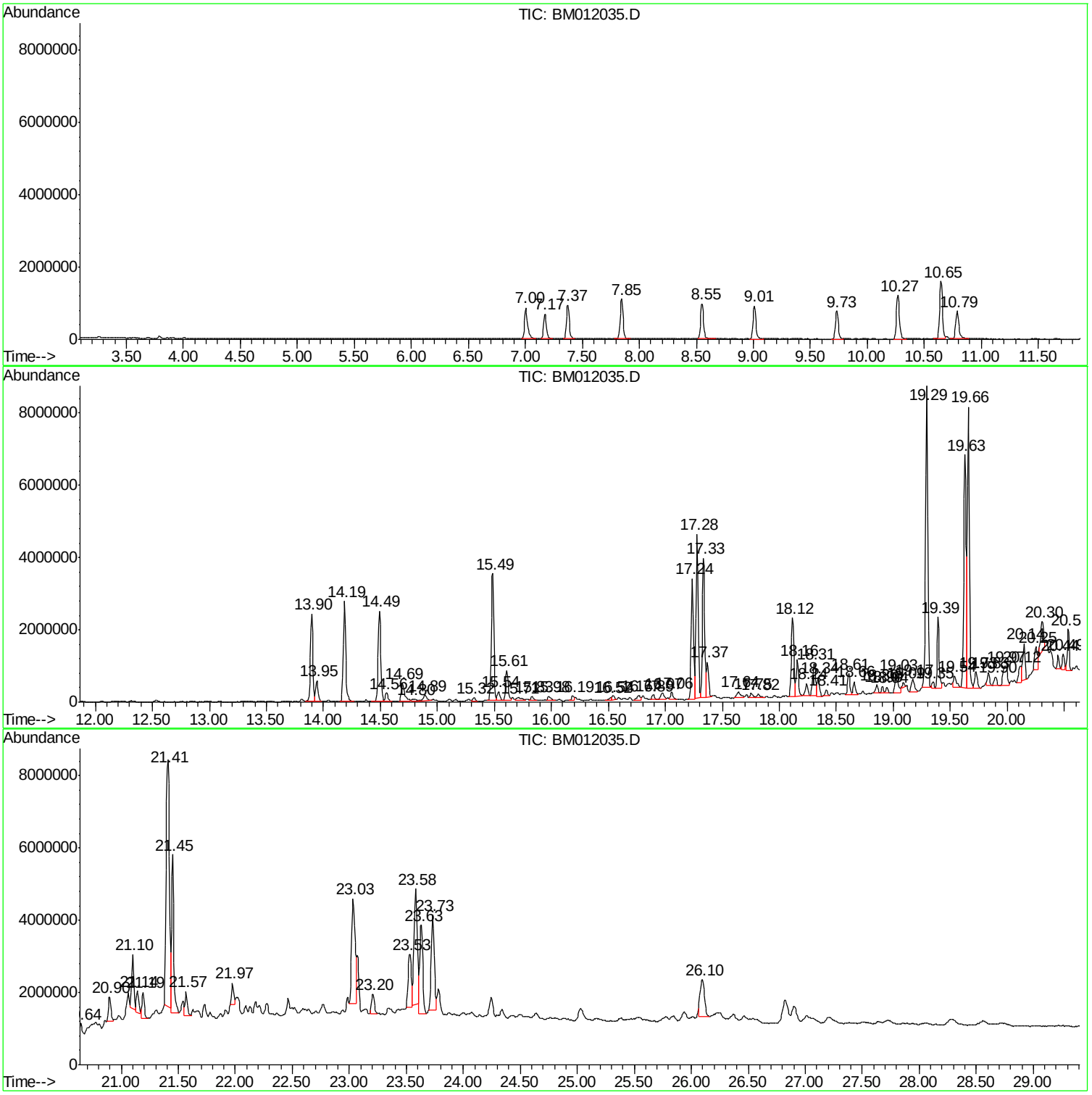
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM100317.M
Quant Title : SVOA CALIBRATION

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



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Operator : SJ/JU
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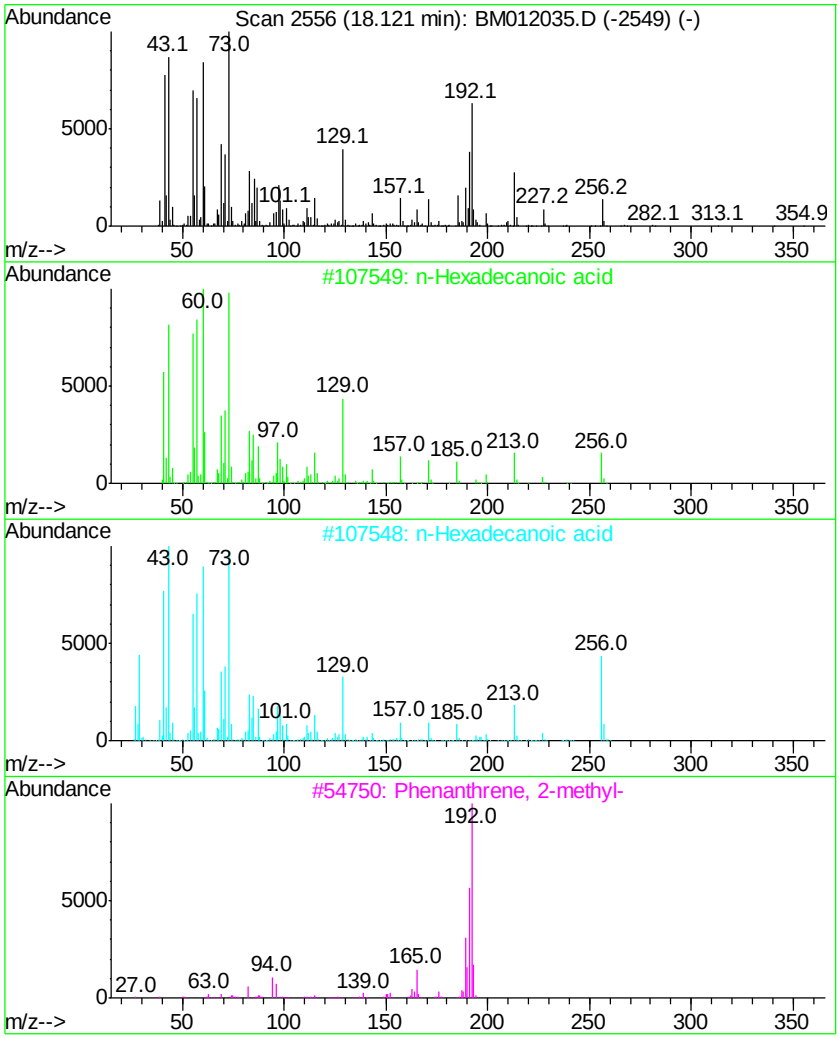
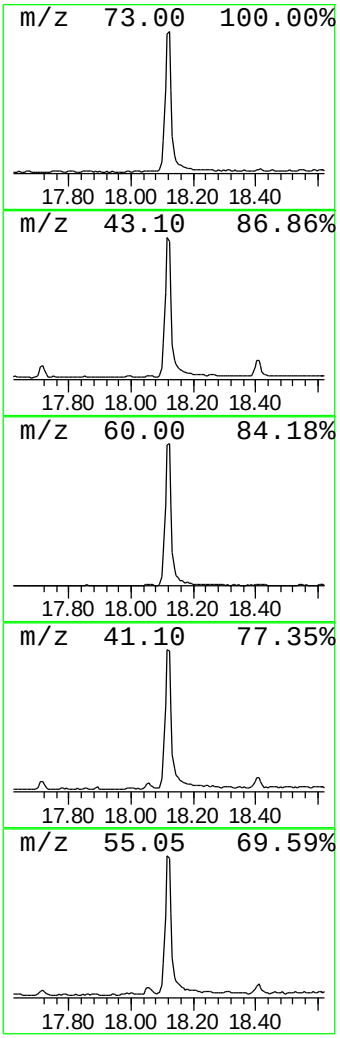
Peak Number 1 n-Hexadecanoic acid Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.12	13.13 ng/ul	3264320	Phenanthrene-d10	17.24

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95
2	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95
3	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	94
4	1H-Cyclopropa[l]phenanthrene,1a,...	192	C15H12	000949-41-7	94
5	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	94





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Sample : I5533-07
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Instrument : BNA_M
ClientSampleId : DUP-1

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM100317.M
Quant Title : SVOA CALIBRATION

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

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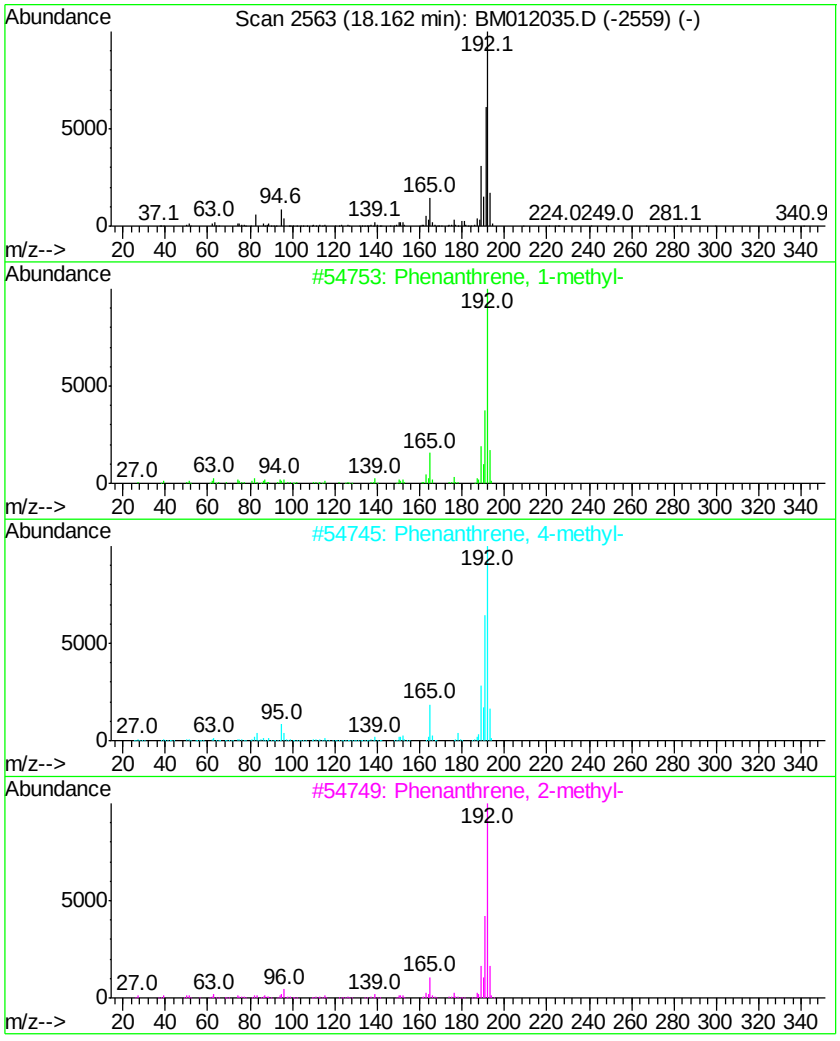
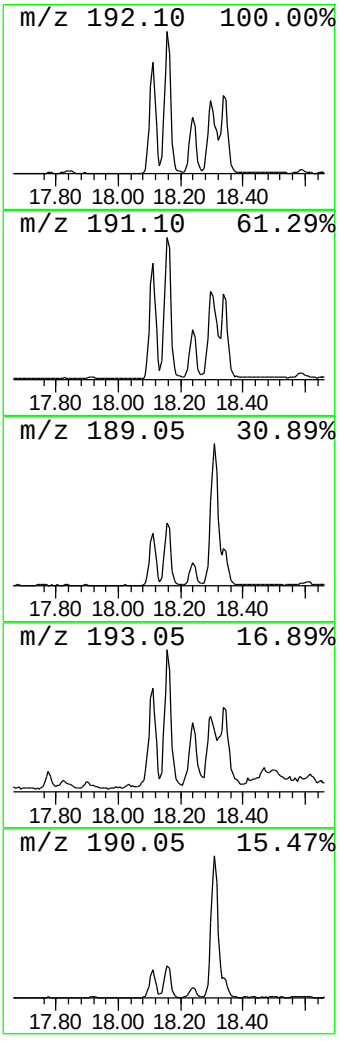
Peak Number 2 Phenanthrene, 1-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.16	5.83 ng/ul	1448780	Phenanthrene-d10	17.24

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	94
2	Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	94
3	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	93
4	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	93
5	Anthracene, 2-methyl-	192	C15H12	000613-12-7	93





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Instrument : BNA_M
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Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM100317.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
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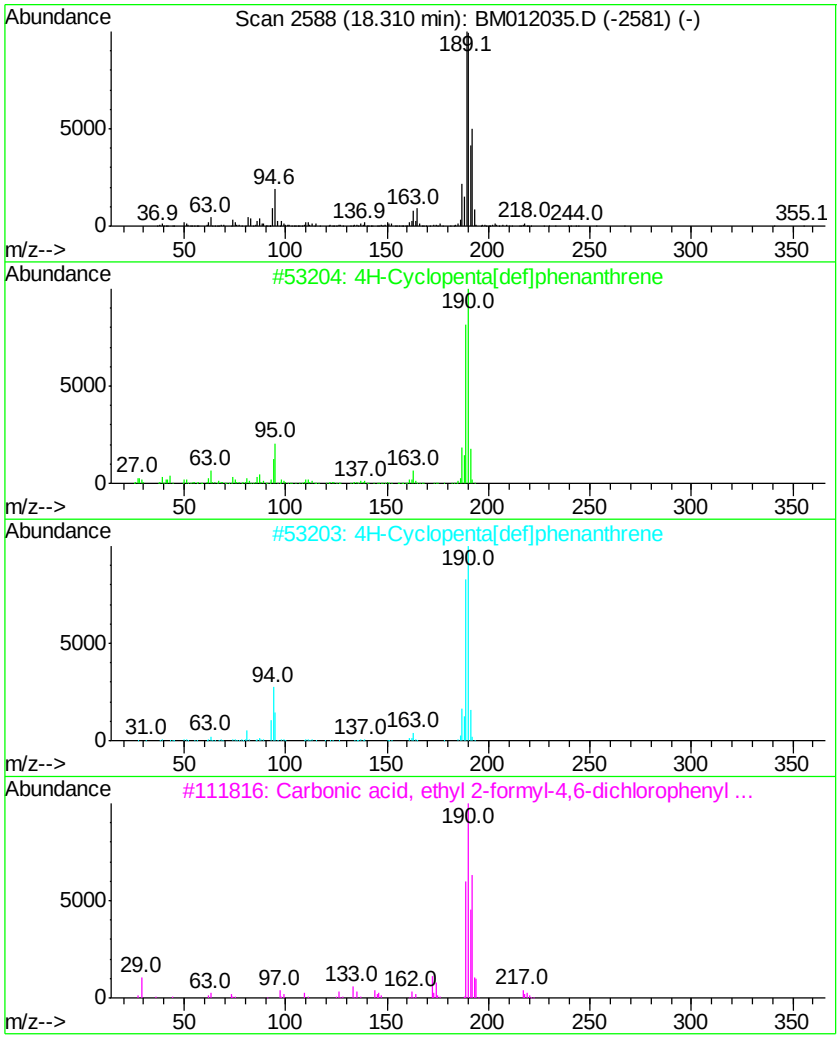
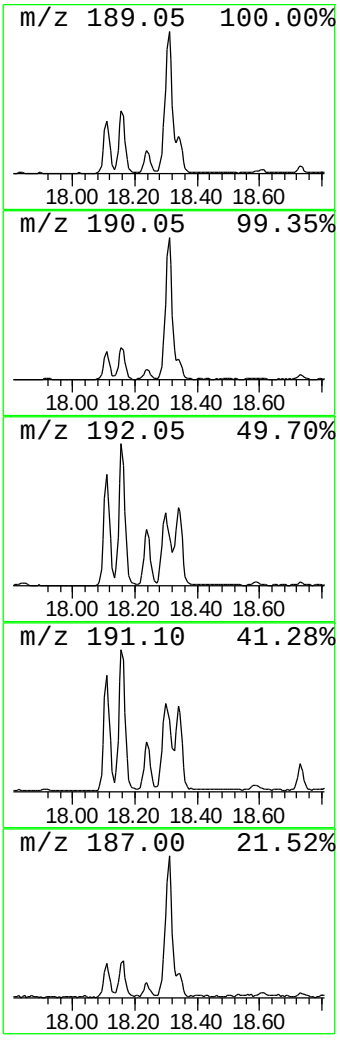
Peak Number 3 4H-Cyclopenta[def]phenanthrene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.31	6.73 ng/ul	1673130	Phenanthrene-d10	17.24

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	64
2	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	62
3	Carbonic acid, ethyl 2-formyl-4,...	262	C10H8Cl2O4	1000331-34-4	50
4	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	43
5	2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	38





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Sample : I5533-07
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : DUP-1

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM100317.M
Quant Title : SVOA CALIBRATION

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

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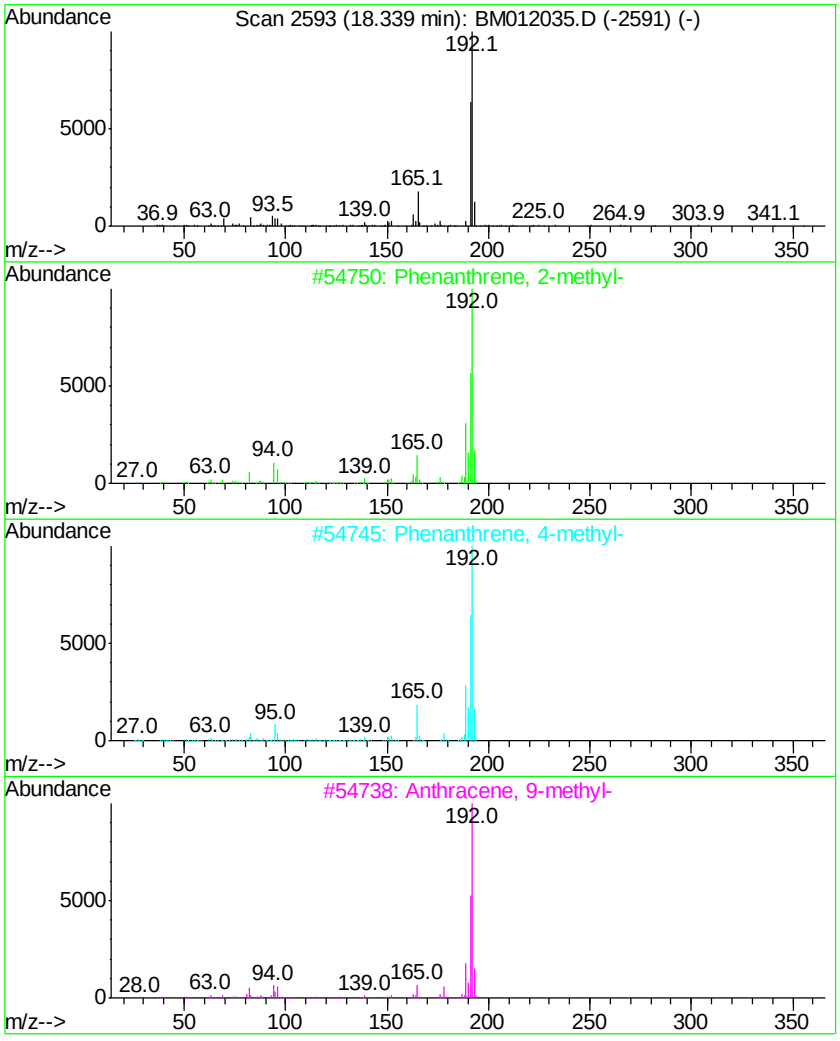
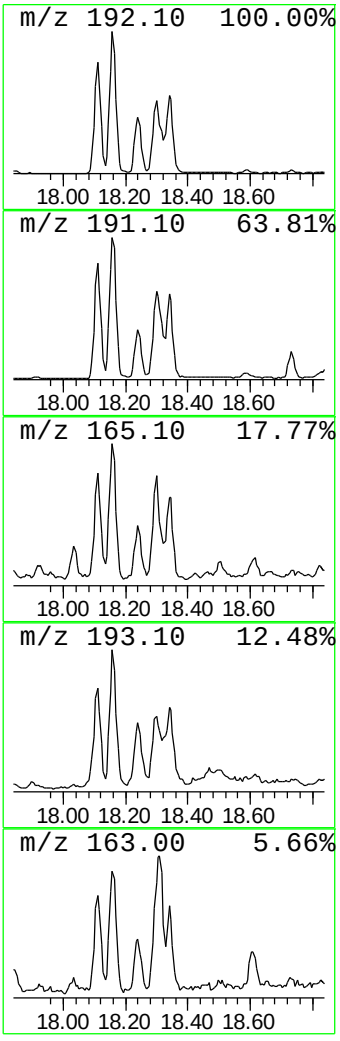
Peak Number 4 Phenanthrene, 2-methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.34	2.70 ng/ul	670456	Phenanthrene-d10	17.24

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	91
2	Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	90
3	Anthracene, 9-methyl-	192	C15H12	000779-02-2	90
4	Anthracene, 1-methyl-	192	C15H12	000610-48-0	87
5	Anthracene, 9-methyl-	192	C15H12	000779-02-2	87





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Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument : BNA_M
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Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM100317.M
Quant Title : SVOA CALIBRATION

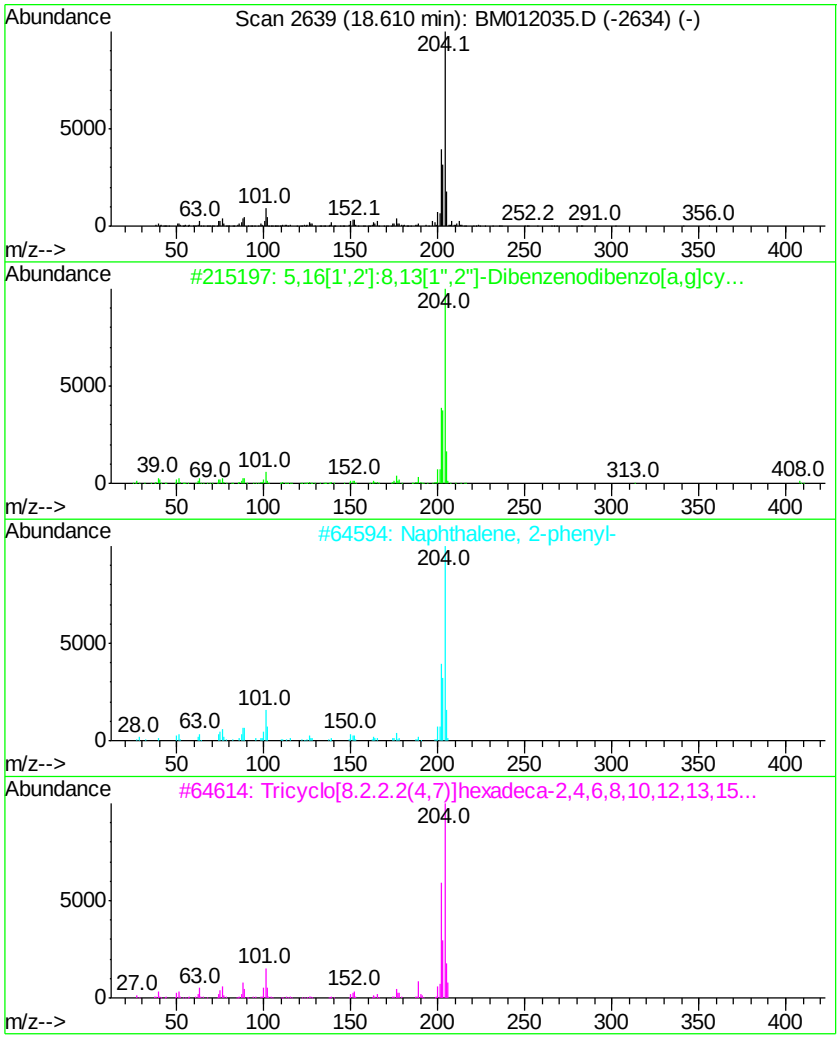
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

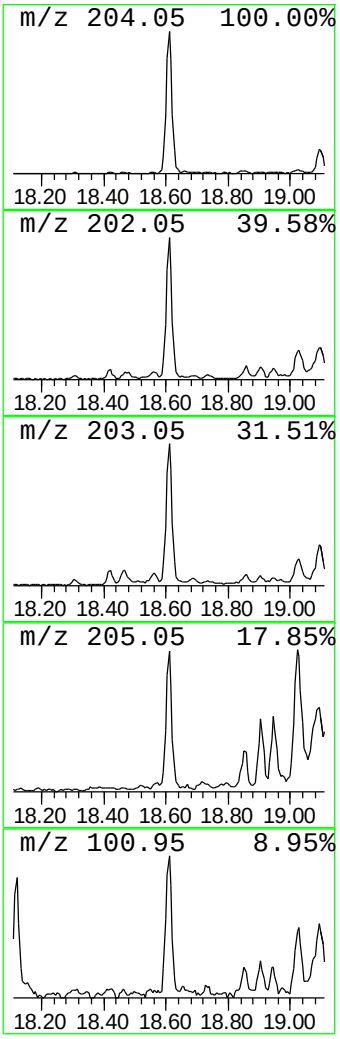
****

Peak Number 5 5,16[1',2']:8,13[1'',2'']-D... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.61	3.10 ng/ul	769781	Phenanthrene-d10	17.24

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5,16[1',2']:8,13[1'',2'']-Dibenz...	408	C32H24	005672-97-9	91
2	Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	86
3	Tricyclo[8.2.2.2(4,7)]hexadeca-2...	204	C16H12	006572-60-7	70
4	Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	70
5	Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	70





Library Search Compound Report

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Data File : BM012035.D
Acq On : 10 Oct 2017 04:40
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Misc :
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Instrument : BNA_M
ClientSampleId : DUP-1

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM100317.M
Quant Title : SVOA CALIBRATION

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

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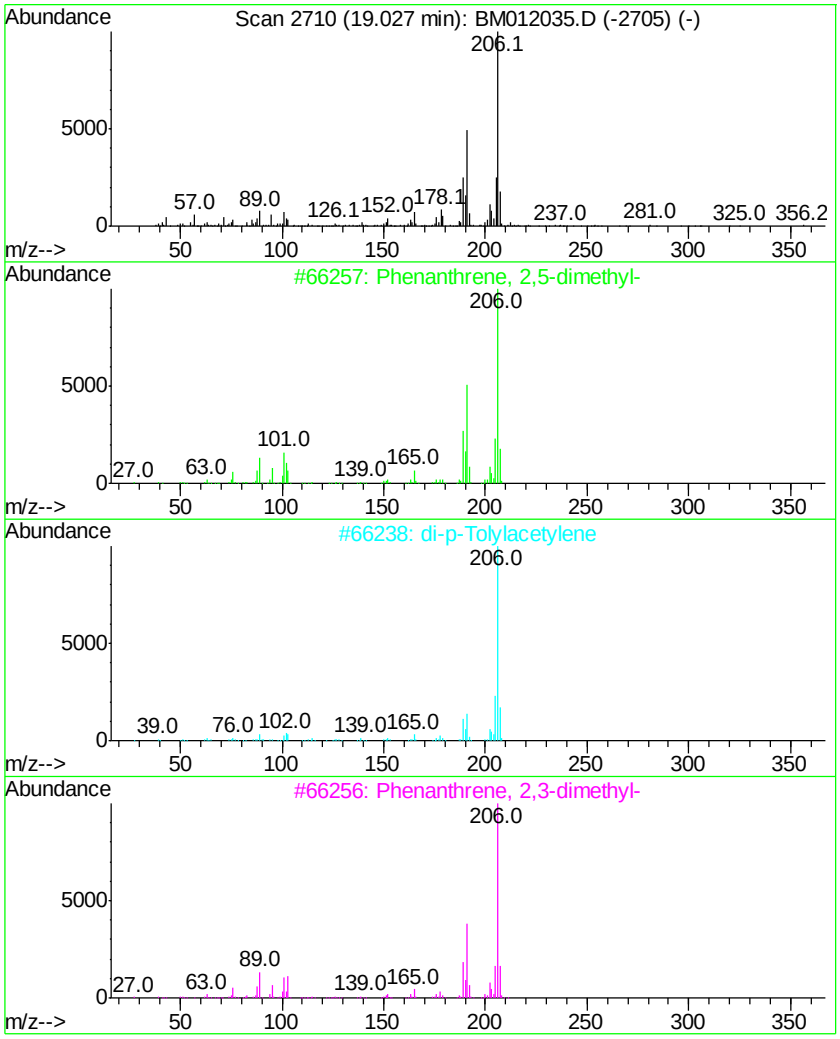
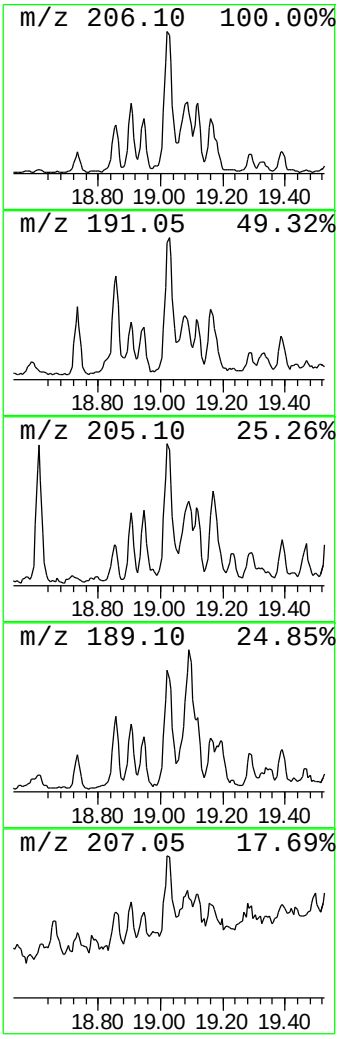
Peak Number 6 Phenanthrene, 2,5-dimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.03	4.19 ng/ul	1041090	Phenanthrene-d10	17.24

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	94
2	di-p-Tolylacetylene	206	C16H14	002789-88-0	91
3	Phenanthrene, 2,3-dimethyl-	206	C16H14	003674-65-5	87
4	9,10-Dimethylanthracene	206	C16H14	000781-43-1	83
5	Naphthalene, 1,2-dihydro-4-phenyl-	206	C16H14	007469-40-1	83





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Data File : BM012035.D
Acq On : 10 Oct 2017 04:40
Operator : SJ/JU
Sample : I5533-07
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : DUP-1

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM100317.M
Quant Title : SVOA CALIBRATION

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

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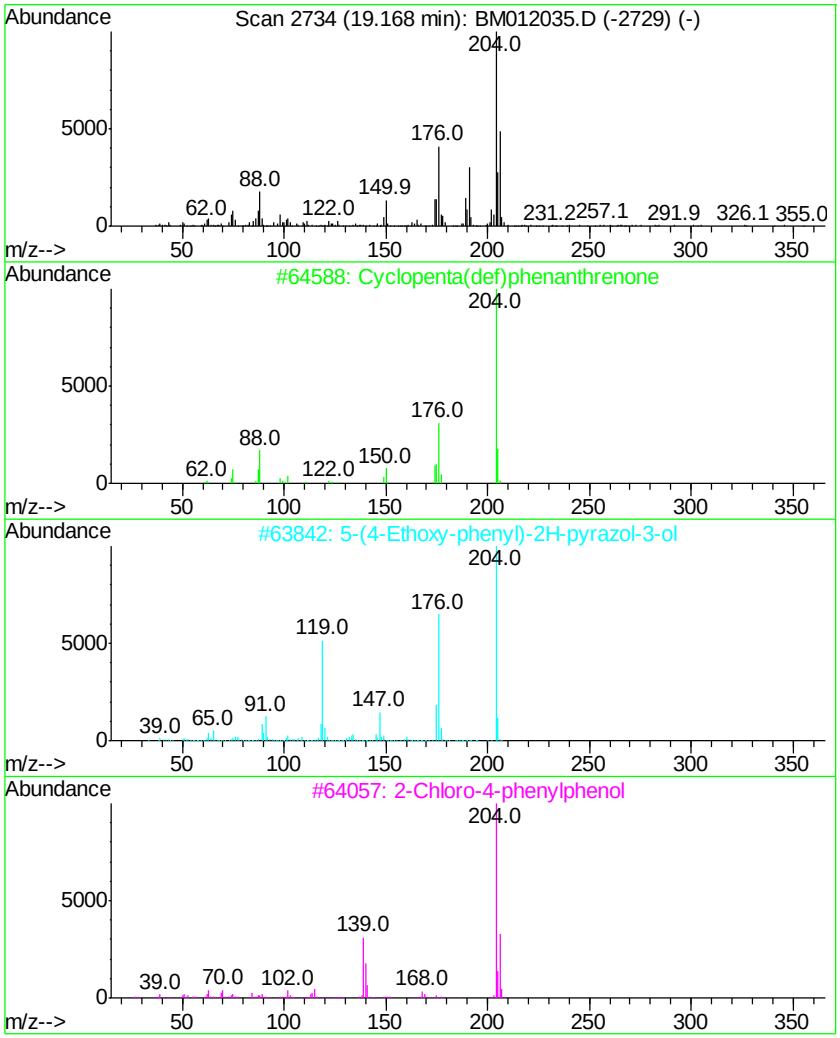
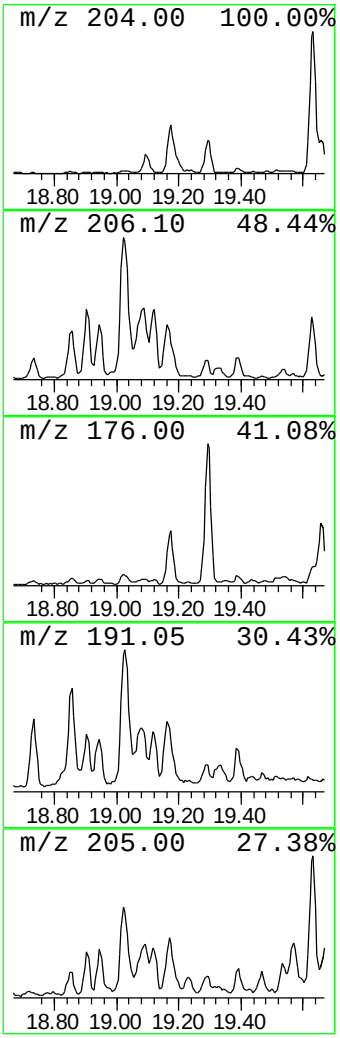
Peak Number 7 Cyclopenta(def)phenanthrenone Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.17	2.78 ng/ul	691914	Phenanthrene-d10	17.24

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopenta(def)phenanthrenone	204	C15H8O	005737-13-3	91
2	5-(4-Ethoxy-phenyl)-2H-pyrazol-3-ol	204	C11H12N2O2	1000278-26-8	43
3	2-Chloro-4-phenylphenol	204	C12H9ClO	000092-04-6	43
4	L-(+)-Rhamnopyranose, tetrakis(t...	452	C18H44O5Si4	1000380-12-9	43
5	1,2,4,8-Tetramethylbicyclo[6.3.0...	204	C15H24	137235-51-9	41





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Operator : SJ/JU
Sample : I5533-07
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : DUP-1

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM100317.M
Quant Title : SVOA CALIBRATION

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

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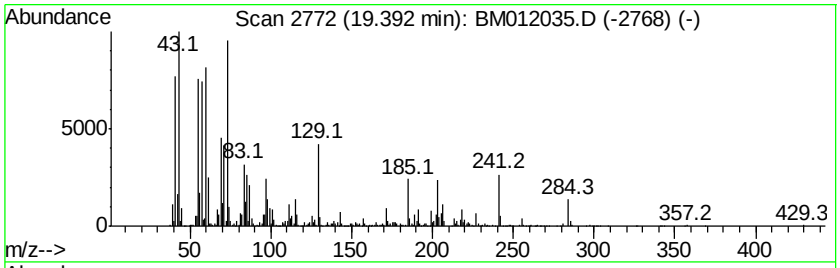
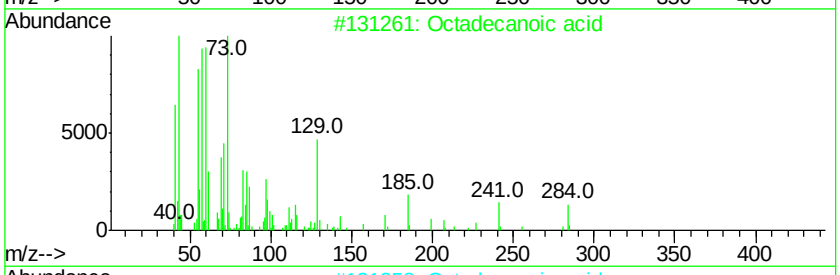
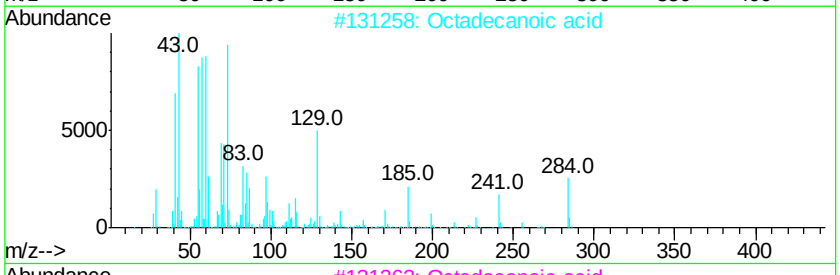
Peak Number 8 Octadecanoic acid Concentration Rank 7

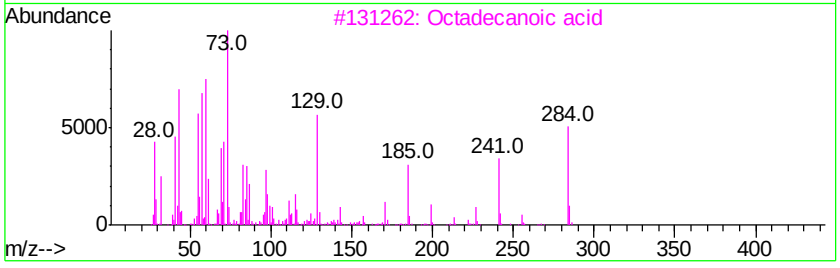
R.T.	EstConc	Area	Relative to ISTD	R.T.
19.39	3.52 ng/ul	2423540	Chrysene-d12	21.42

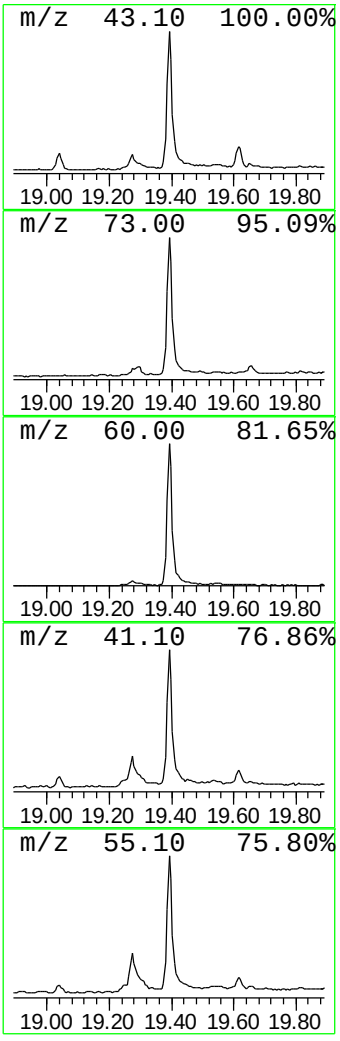
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanoic acid	284	C18H36O2	000057-11-4	99
2	Octadecanoic acid	284	C18H36O2	000057-11-4	95
3	Octadecanoic acid	284	C18H36O2	000057-11-4	95
4	Pentadecanoic acid	242	C15H30O2	001002-84-2	86
5	Pentadecanoic acid	242	C15H30O2	001002-84-2	78











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Operator : SJ/JU
Sample : I5533-07
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : DUP-1

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM100317.M
Quant Title : SVOA CALIBRATION

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

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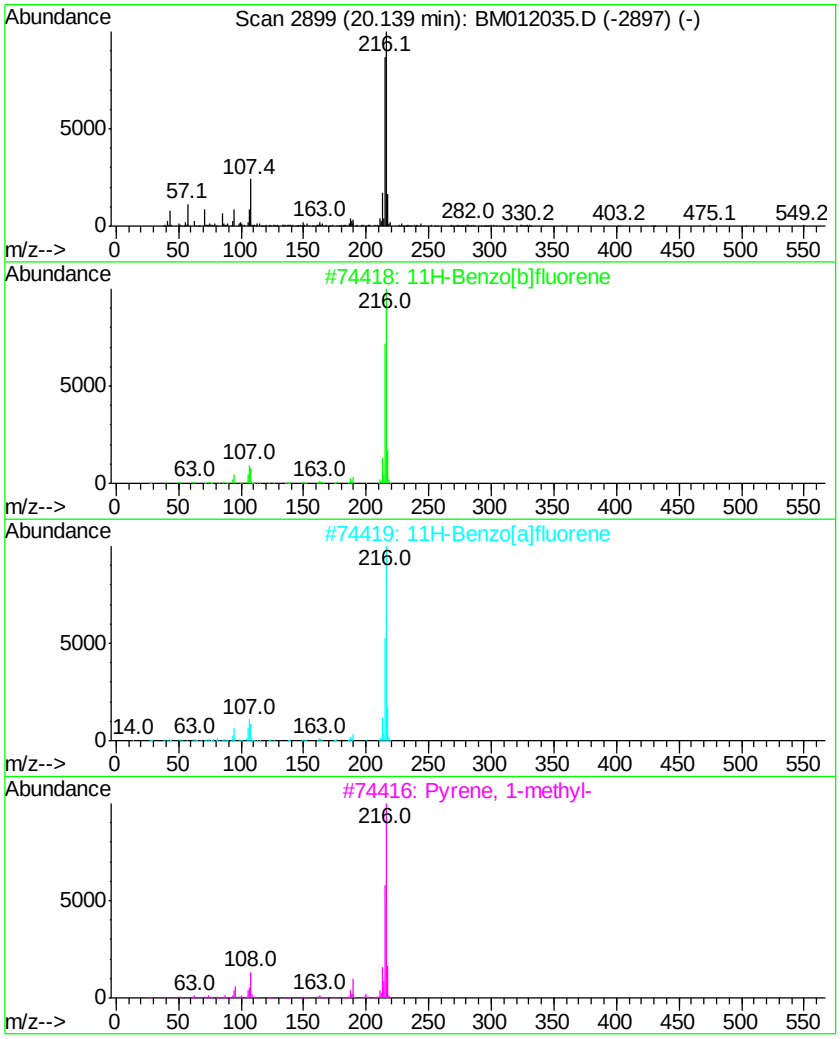
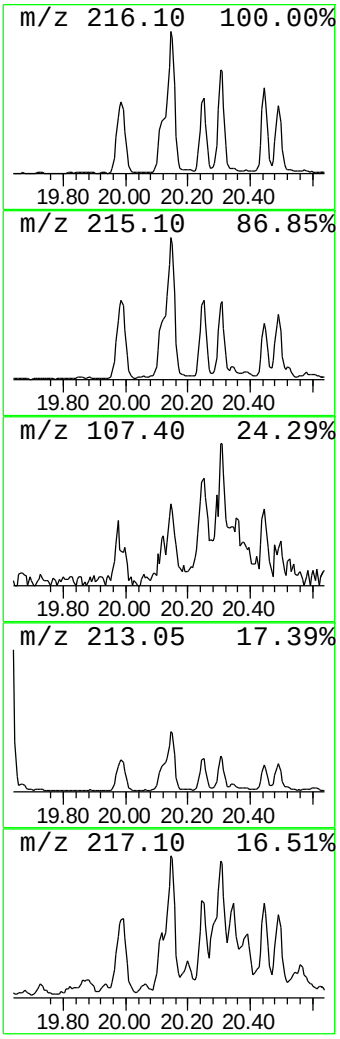
Peak Number 9 11H-Benzo[b]fluorene Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.14	2.06 ng/ul	1415560	Chrysene-d12	21.42

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	95
2	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	90
3	Pyrene, 1-methyl-	216	C17H12	002381-21-7	87
4	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	83
5	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	83





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Misc :
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Instrument : BNA_M
ClientSampleId : DUP-1

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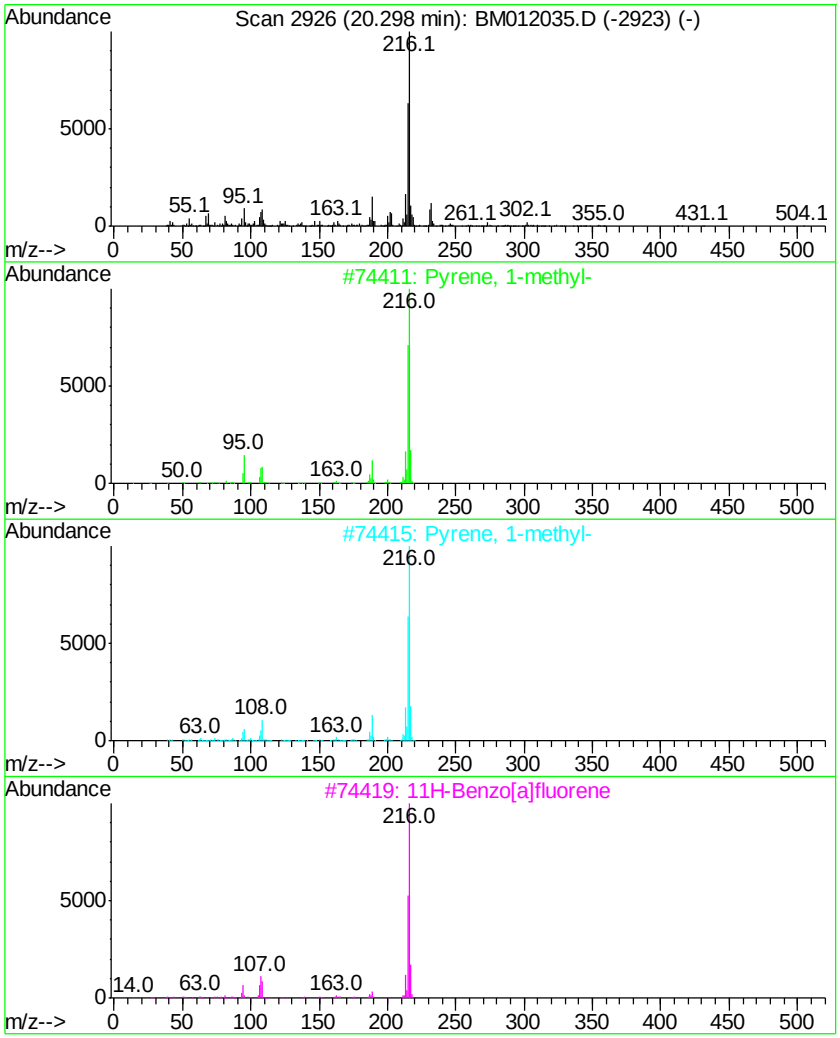
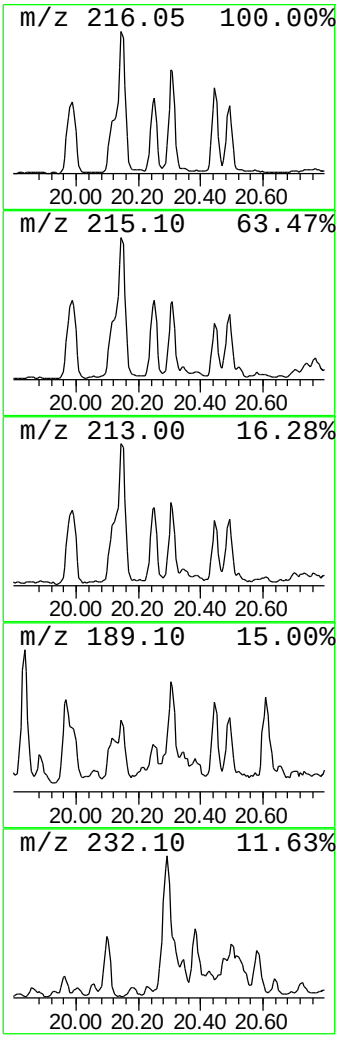
Peak Number 10 Pyrene, 1-methyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.30	2.08 ng/ul	1433360	Chrysene-d12	21.42

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pyrene, 1-methyl-	216	C17H12	002381-21-7	97
2	Pyrene, 1-methyl-	216	C17H12	002381-21-7	96
3	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	94
4	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	93
5	Pyrene, 2-methyl-	216	C17H12	003442-78-2	93





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Operator : SJ/JU
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Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : DUP-1

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM100317.M
Quant Title : SVOA CALIBRATION

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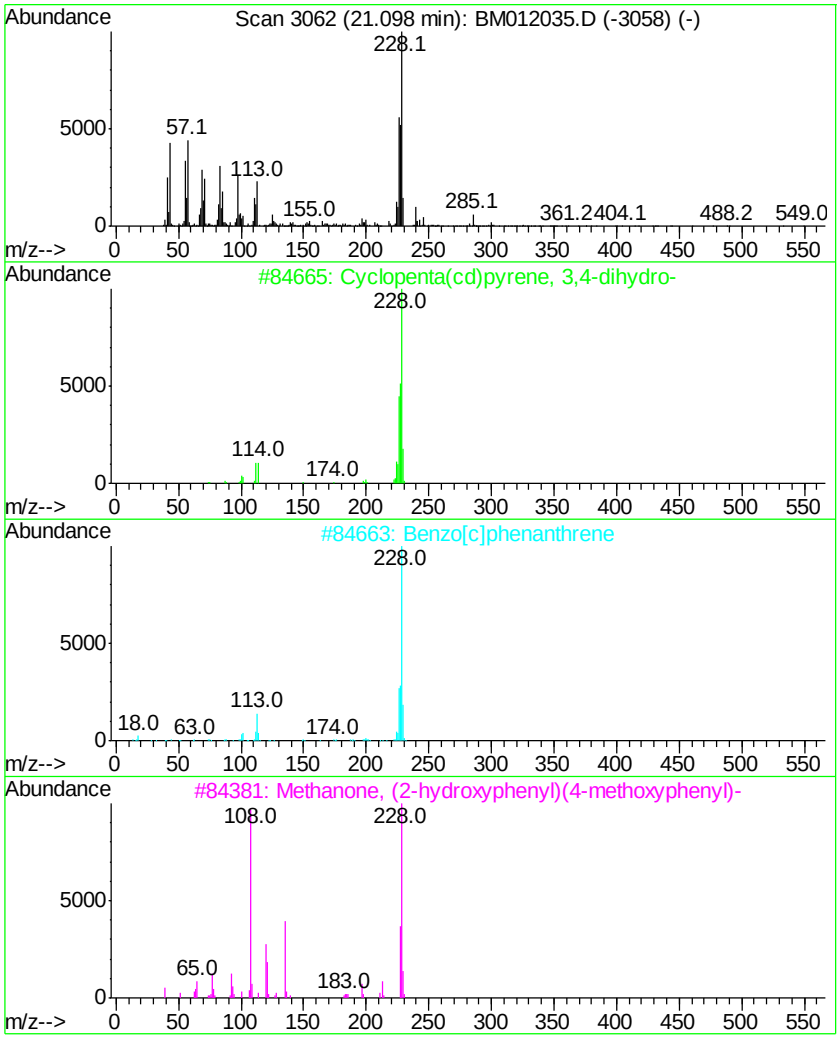
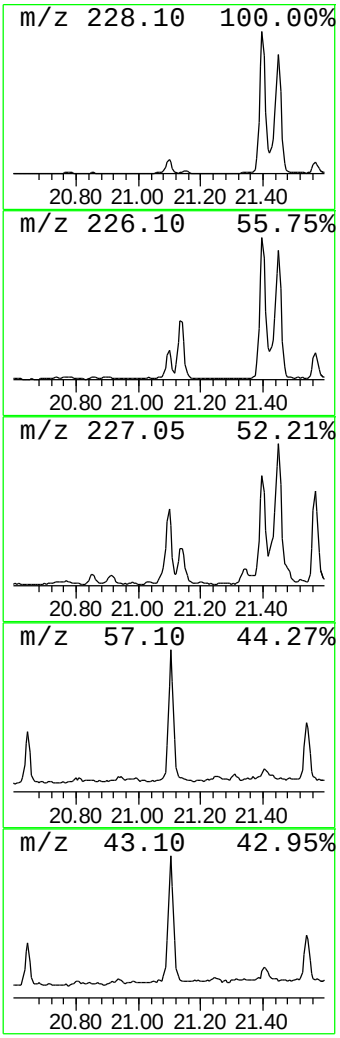
Peak Number 11 Cyclopenta(cd)pyrene, 3,4-d... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.10	2.87 ng/ul	1976860	Chrysene-d12	21.42

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopenta(cd)pyrene, 3,4-dihydro-	228	C18H12	025732-74-5	50
2	Benzo[c]phenanthrene	228	C18H12	000195-19-7	49
3	Methanone, (2-hydroxyphenyl)(4-m...	228	C14H12O3	018733-07-8	46
4	Purine-6(1H)-thione, 3,7-dimethy...	228	C7H8N4Se	023663-58-3	38
5	Triphenylene	228	C18H12	000217-59-4	38





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Misc :
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Instrument : BNA_M
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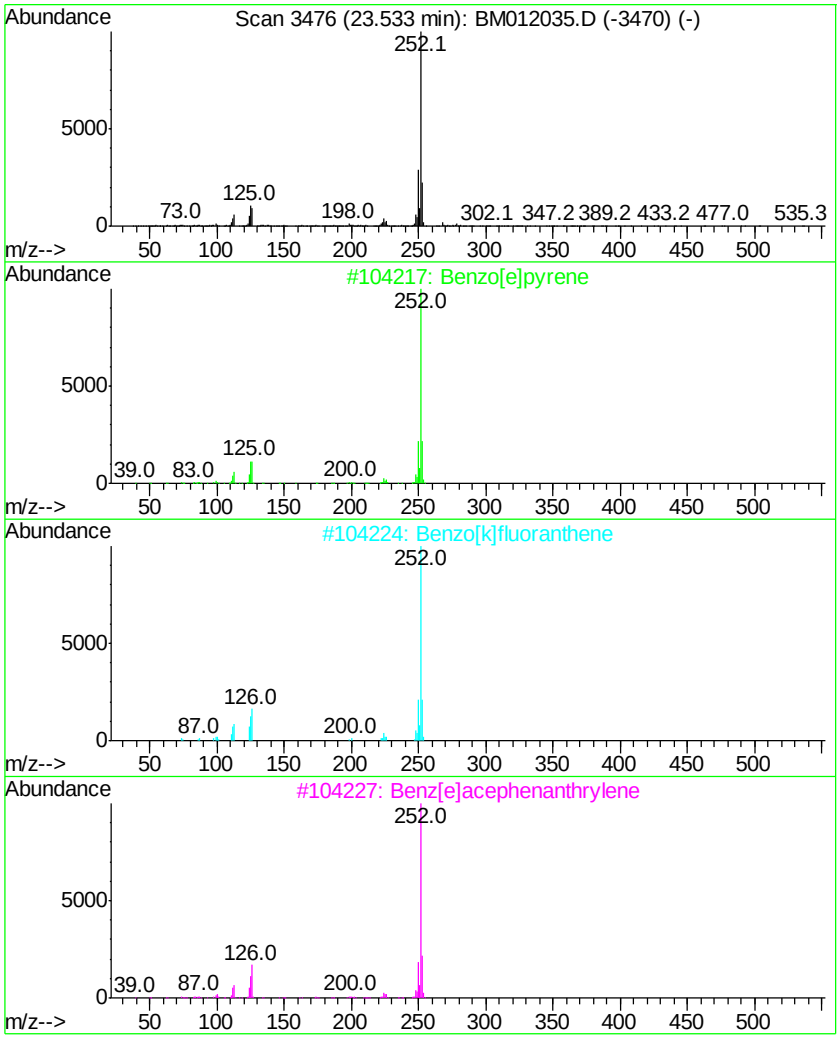
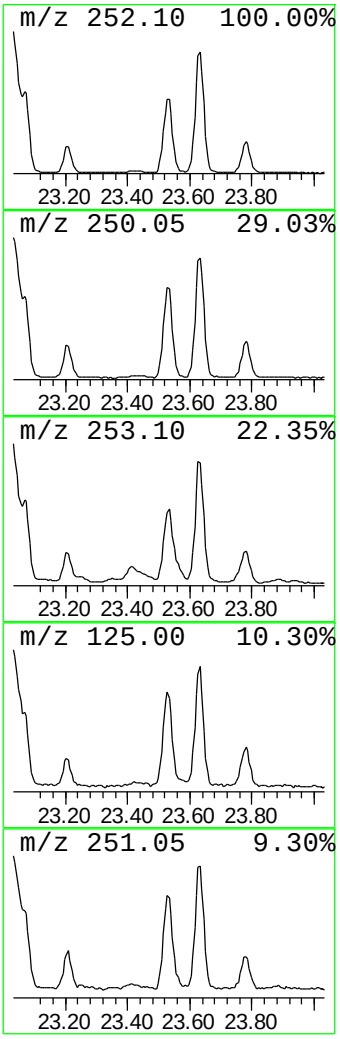
Peak Number 13 Benzo[e]pyrene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.53	11.45 ng/ul	2860480	Perylene-d12	23.73

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzo[e]pyrene	252	C20H12	000192-97-2	98
2	Benzo[k]fluoranthene	252	C20H12	000207-08-9	95
3	Benz[e]acephenanthrylene	252	C20H12	000205-99-2	93
4	Benzo[a]pyrene	252	C20H12	000050-32-8	93
5	Perylene	252	C20H12	000198-55-0	90





Tentatively Identified Compound (LSC) summary

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM100917\
Data File : BM012035.D
Acq On : 10 Oct 2017 04:40
Operator : SJ/JU
Sample : I5533-07
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : DUP-1

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM100317.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
n-Hexadecanoic acid	18.12	13.1	ng/ul	3264320	4	17.24	4971250	20.0
Phenanthrene, 1-m...	18.16	5.8	ng/ul	1448780	4	17.24	4971250	20.0
4H-Cyclopenta[def...	18.31	6.7	ng/ul	1673130	4	17.24	4971250	20.0
Phenanthrene, 2-m...	18.34	2.7	ng/ul	670456	4	17.24	4971250	20.0
5,16[1',2']:8,13[...	18.61	3.1	ng/ul	769781	4	17.24	4971250	20.0
Phenanthrene, 2,5...	19.03	4.2	ng/ul	1041090	4	17.24	4971250	20.0
Cyclopenta(def)ph...	19.17	2.8	ng/ul	691914	4	17.24	4971250	20.0
Octadecanoic acid	19.39	3.5	ng/ul	2423540	5	21.42	13766600	20.0
11H-Benzo[b]fluorene	20.14	2.1	ng/ul	1415560	5	21.42	13766600	20.0
Pyrene, 1-methyl-	20.30	2.1	ng/ul	1433360	5	21.42	13766600	20.0
Cyclopenta(cd)pyr...	21.10	2.9	ng/ul	1976860	5	21.42	13766600	20.0
Benzo[e]pyrene	23.53	11.4	ng/ul	2860480	6	23.73	4997210	20.0