

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM062317\
 Data File : BM010708.D
 Acq On : 24 Jun 2017 02:02
 Operator : SJ/JU
 Sample : I3625-18
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 F5A00

Integration Parameters: LSCINT.P

Integrator: RTE
 Smoothing : OFF
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0

Filtering: 5
 Min Area: 1 % of largest Peak
 Max Peaks: 100
 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	6.645	632	639	649	rBB	326422	524633	14.99%	1.090%
2	6.810	662	667	676	rVB	222417	327416	9.36%	0.680%
3	6.998	692	699	707	rBB	327605	490532	14.02%	1.019%
4	7.457	770	777	784	rBV	365287	565354	16.16%	1.174%
5	8.163	891	897	905	rBV	439935	666589	19.05%	1.384%
6	8.604	965	972	980	rBV	332034	522989	14.95%	1.086%
7	9.316	1087	1093	1100	rVB	314064	500391	14.30%	1.039%
8	9.845	1175	1183	1191	rBB	504506	792864	22.66%	1.647%
9	10.222	1240	1247	1252	rBV	571183	905994	25.89%	1.882%
10	10.269	1252	1255	1264	rVB	38373	58252	1.66%	0.121%
11	10.363	1264	1271	1279	rBV	434292	686193	19.61%	1.425%
12	13.533	1804	1810	1814	rBV	1004293	1305324	37.31%	2.711%
13	13.580	1814	1818	1824	rVB	304071	396418	11.33%	0.823%
14	13.798	1849	1855	1865	rVB	1010214	1454237	41.56%	3.020%
15	14.110	1902	1908	1915	rBV3	892938	1366417	39.05%	2.838%
16	14.174	1915	1919	1924	rVV	46666	60761	1.74%	0.126%
17	14.321	1938	1944	1953	rBV	434937	587176	16.78%	1.220%
18	14.515	1972	1977	1981	rVB2	34768	44012	1.26%	0.091%
19	15.115	2072	2079	2084	rBV	1251559	1808320	51.68%	3.756%
20	15.168	2084	2088	2094	rVB	68809	92105	2.63%	0.191%
21	15.239	2094	2100	2106	rBV	421030	562047	16.06%	1.167%
22	16.521	2313	2318	2325	rVB	53721	88378	2.53%	0.184%
23	16.621	2328	2335	2339	rBV5	34748	66693	1.91%	0.139%
24	16.680	2343	2345	2350	rVB	38423	45586	1.30%	0.095%
25	16.868	2371	2377	2380	rBV	1220833	1695802	48.46%	3.522%
26	16.909	2380	2384	2388	rVV	1140271	1511922	43.21%	3.140%
27	16.968	2388	2394	2398	rVV	1386036	2031417	58.06%	4.219%
28	16.998	2398	2399	2405	rVV	181911	158169	4.52%	0.329%
29	17.274	2442	2446	2450	rVB	86019	111094	3.17%	0.231%
30	17.392	2463	2466	2471	rBV3	32809	45374	1.30%	0.094%
31	17.445	2471	2475	2481	rVB6	29716	44258	1.26%	0.092%
32	17.745	2520	2526	2529	rVV	185033	262482	7.50%	0.545%
33	17.792	2529	2534	2541	rVV	388058	545767	15.60%	1.134%
34	17.868	2542	2547	2552	rVV2	74427	108100	3.09%	0.225%

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35	17.933	2552	2558	2562	rVV2	227303	408515	11.68%	0.848%
36	17.974	2562	2565	2572	rVB	137716	184786	5.28%	0.384%
37	18.251	2608	2612	2615	rBV	157116	200510	5.73%	0.416%
38	18.286	2615	2618	2625	rVB	100598	146203	4.18%	0.304%
39	18.503	2651	2655	2659	rVB3	51783	66835	1.91%	0.139%
40	18.551	2659	2663	2667	rVB	42802	52563	1.50%	0.109%
41	18.592	2667	2670	2675	rVB2	44735	53572	1.53%	0.111%
42	18.674	2675	2684	2689	rBV2	181632	297516	8.50%	0.618%
43	18.733	2689	2694	2697	rVV3	74217	121845	3.48%	0.253%
44	18.768	2697	2700	2703	rVB	40076	48404	1.38%	0.101%
45	18.815	2703	2708	2716	rVB2	106559	186840	5.34%	0.388%
46	18.939	2720	2729	2738	rVB	2055191	2656546	75.92%	5.517%
47	19.015	2738	2742	2744	rBV	45607	72534	2.07%	0.151%
48	19.039	2744	2746	2750	rVV5	45931	66880	1.91%	0.139%
49	19.086	2750	2754	2758	rVV2	137642	168526	4.82%	0.350%
50	19.145	2758	2764	2768	rVV5	47688	103706	2.96%	0.215%
51	19.180	2768	2770	2775	rVB	56915	73001	2.09%	0.152%
52	19.274	2781	2786	2789	rBV2	1947692	2892347	82.66%	6.007%
53	19.303	2789	2791	2799	rVV	1648737	2010990	57.47%	4.177%
54	19.380	2799	2804	2811	rVB3	92260	169509	4.84%	0.352%
55	19.486	2816	2822	2827	rBV	104087	152205	4.35%	0.316%
56	19.545	2827	2832	2838	rVB2	72106	128973	3.69%	0.268%
57	19.639	2838	2848	2854	rBV3	157952	401001	11.46%	0.833%
58	19.774	2866	2871	2873	rBV2	106385	188486	5.39%	0.391%
59	19.803	2873	2876	2881	rVB	209753	315903	9.03%	0.656%
60	19.862	2881	2886	2889	rBV6	24514	51454	1.47%	0.107%
61	19.909	2890	2894	2897	rBV2	81245	91212	2.61%	0.189%
62	19.962	2897	2903	2907	rBV2	213482	333379	9.53%	0.692%
63	20.003	2907	2910	2914	rVB2	58870	71881	2.05%	0.149%
64	20.050	2914	2918	2922	rVB3	42897	52757	1.51%	0.110%
65	20.103	2923	2927	2931	rBV	117698	153956	4.40%	0.320%
66	20.150	2931	2935	2941	rBV	138259	177127	5.06%	0.368%
67	20.262	2949	2954	2961	rVB10	27751	63865	1.83%	0.133%
68	20.368	2968	2972	2973	rBV3	47149	54484	1.56%	0.113%
69	20.403	2974	2978	2981	rVV5	56079	100647	2.88%	0.209%
70	20.450	2983	2986	2988	rVV	42667	52352	1.50%	0.109%
71	20.556	2994	3004	3012	rBV	262156	422196	12.07%	0.877%

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72	20.721	3026	3032	3036	rBV2	201745	370325	10.58%	0.769%
73	20.762	3036	3039	3043	rVV	175821	238190	6.81%	0.495%
74	20.815	3043	3048	3052	rVV3	235096	442559	12.65%	0.919%
75	20.856	3052	3055	3062	rVB	226781	286493	8.19%	0.595%
76	20.962	3067	3073	3079	rBV	69746	136563	3.90%	0.284%
77	21.033	3082	3085	3086	rBV2	59862	61848	1.77%	0.128%
78	21.080	3086	3093	3097	rVV3	1831328	3499046	100.00%	7.267%
79	21.121	3097	3100	3111	rVB	915247	1216001	34.75%	2.526%
80	21.233	3117	3119	3122	rBV	54013	65979	1.89%	0.137%
81	21.392	3141	3146	3151	rBV3	129386	236184	6.75%	0.491%
82	21.439	3151	3154	3157	rVB4	53318	72962	2.09%	0.152%
83	21.568	3173	3176	3178	rBV2	53552	57876	1.65%	0.120%
84	21.621	3179	3185	3190	rVV	189661	291490	8.33%	0.605%
85	21.674	3190	3194	3199	rVV2	116280	207127	5.92%	0.430%
86	21.809	3213	3217	3219	rBV	89981	114605	3.28%	0.238%
87	21.839	3219	3222	3228	rVB4	75156	99670	2.85%	0.207%
88	22.127	3268	3271	3276	rVB5	93002	124755	3.57%	0.259%
89	22.356	3302	3310	3316	rBV4	54719	150443	4.30%	0.312%
90	22.591	3344	3350	3361	rBV	621860	1736762	49.64%	3.607%
91	22.750	3373	3377	3383	rVB	78910	120979	3.46%	0.251%
92	22.874	3395	3398	3405	rBV5	40177	83640	2.39%	0.174%
93	23.038	3421	3426	3430	rBV	296968	543260	15.53%	1.128%
94	23.091	3430	3435	3439	rVV	910019	1614725	46.15%	3.354%
95	23.133	3439	3442	3450	rVB	491196	762716	21.80%	1.584%
96	23.221	3451	3457	3462	rBV	701213	1275645	36.46%	2.649%
97	23.268	3463	3465	3472	rVB2	132052	197430	5.64%	0.410%
98	23.750	3543	3547	3552	rVB5	72453	129835	3.71%	0.270%
99	25.332	3809	3816	3826	rVB3	197332	505256	14.44%	1.049%
100	25.979	3920	3926	3935	rVB2	104349	279306	7.98%	0.580%

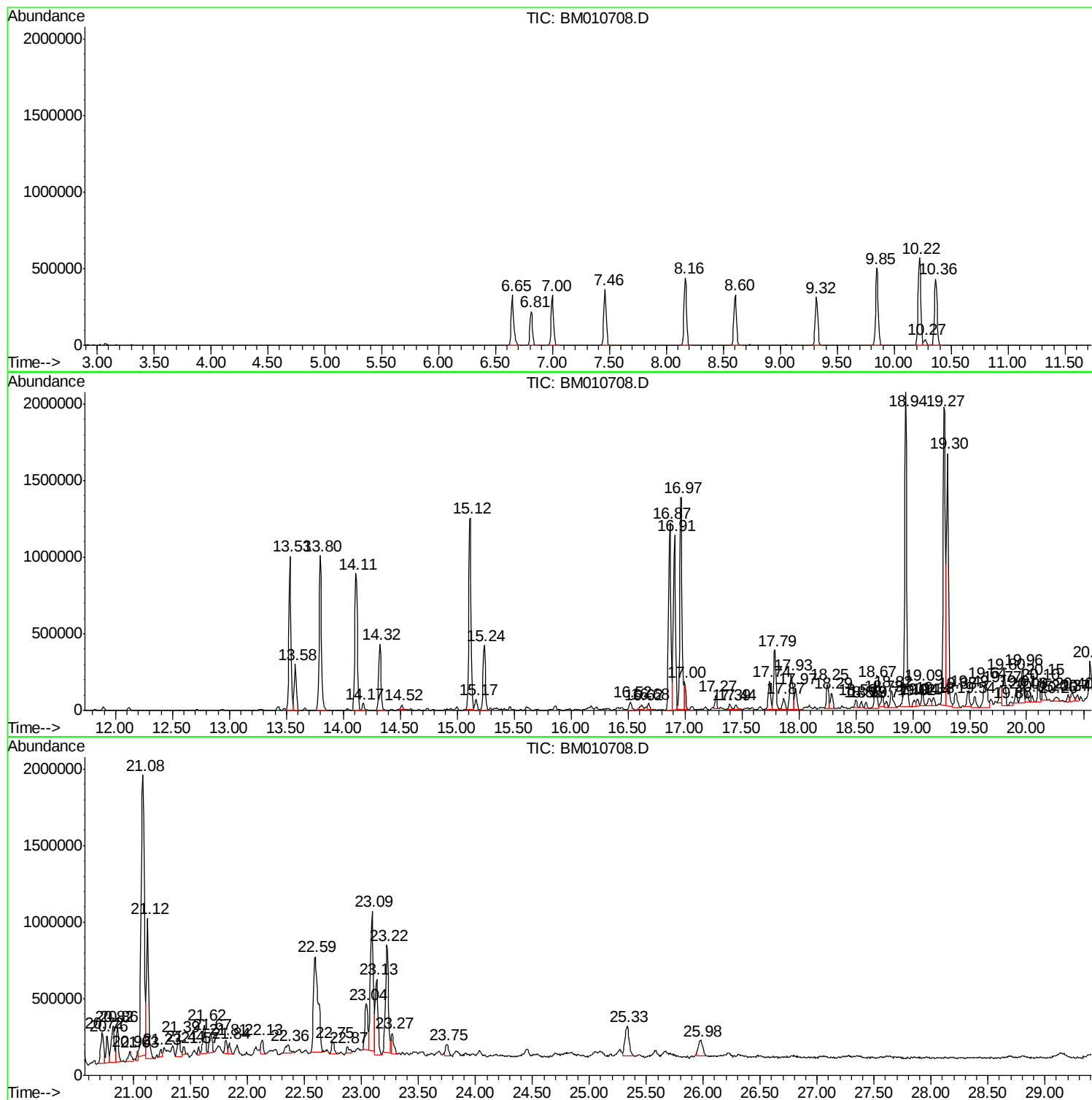
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Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM062017.M
Quant Title : SVOA CALIBRATION

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



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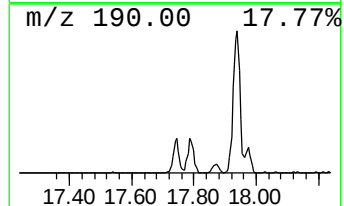
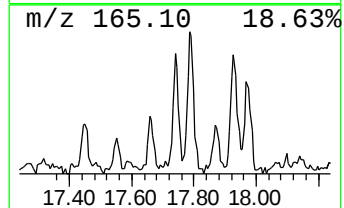
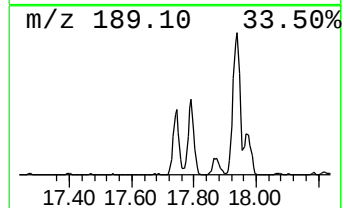
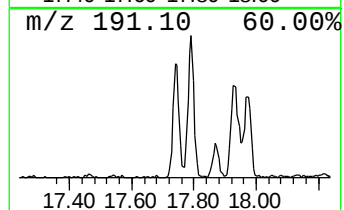
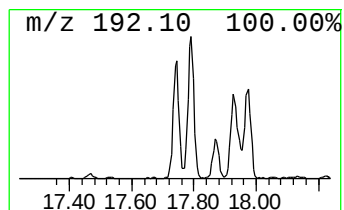
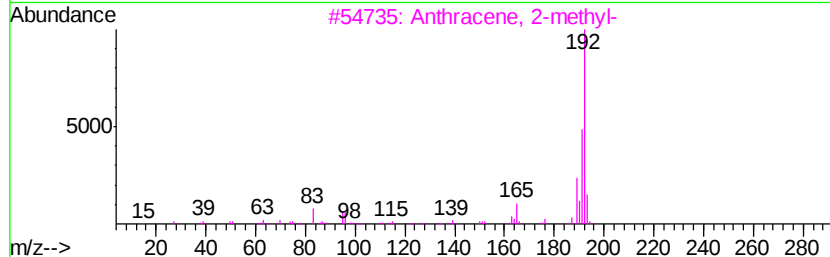
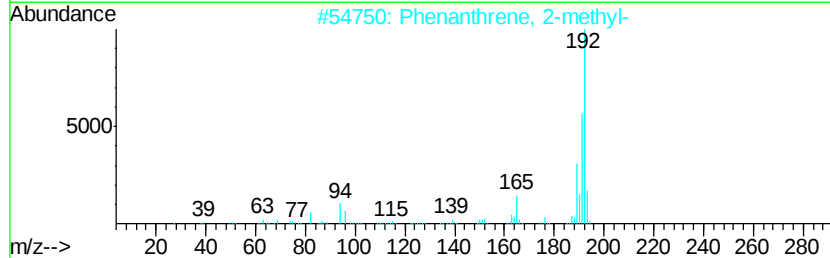
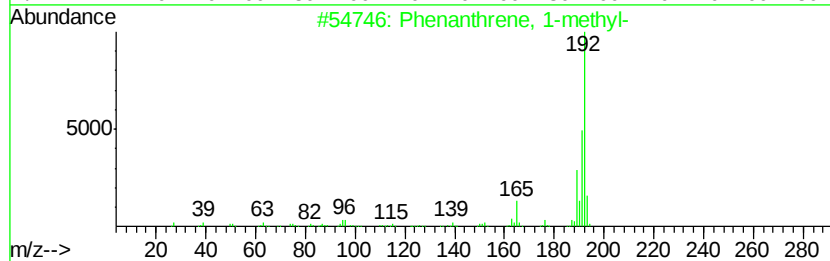
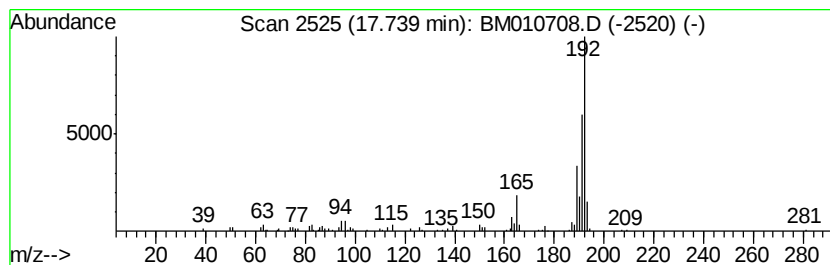
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Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM062017.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Phenanthrene, 1-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
17.74	3.10 ng/ul	262482	Phenanthrene-d10	16.87	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	96
2	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	95
3	Anthracene, 2-methyl-	192	C15H12	000613-12-7	94
4	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	94
5	Anthracene, 2-methyl-	192	C15H12	000613-12-7	94



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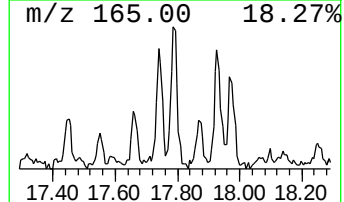
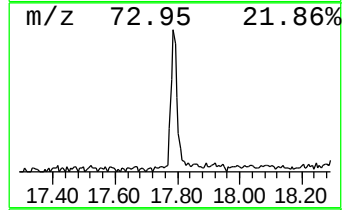
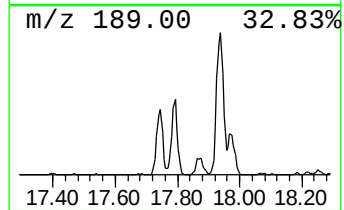
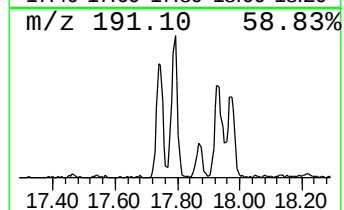
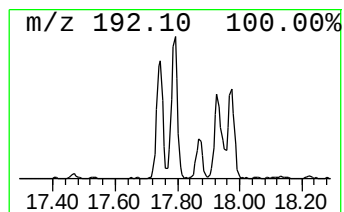
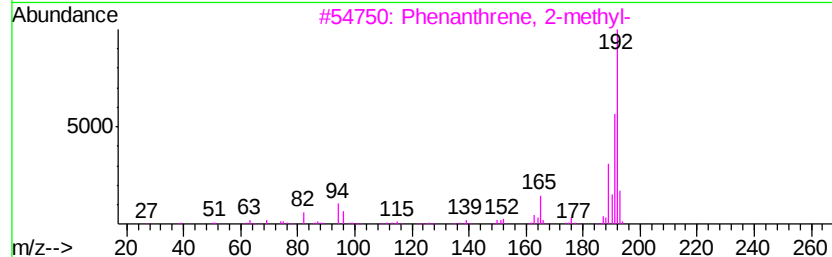
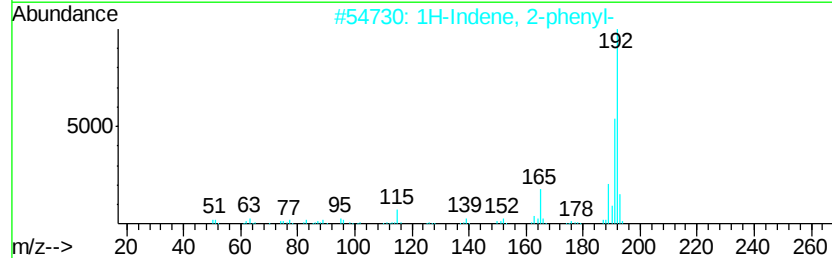
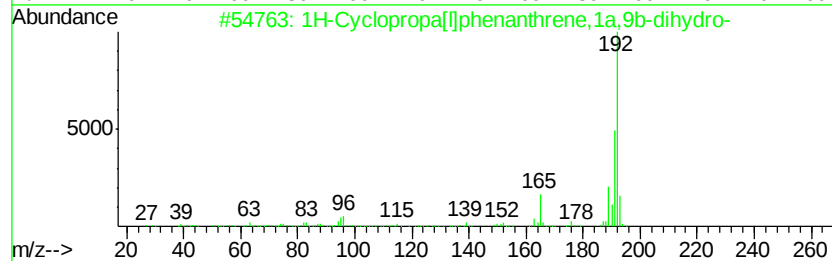
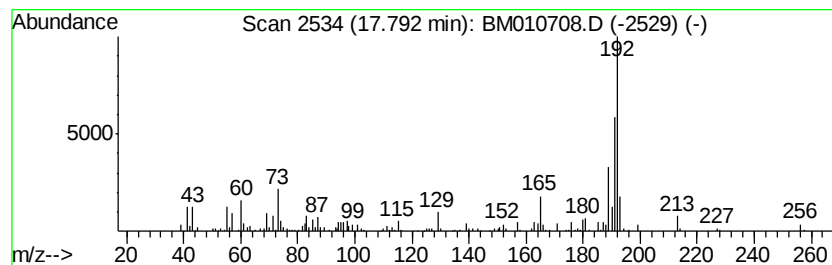
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 1H-Cyclopropa[l]phenanthren... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.79	6.44 ng/ul	545767	Phenanthrene-d10	16.87

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		1H-Cyclopropa[l]phenanthrene, 1a, ...	192	C15H12	000949-41-7	97
2		1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	96
3		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	95
4		Anthracene, 2-methyl-	192	C15H12	000613-12-7	95
5		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	95



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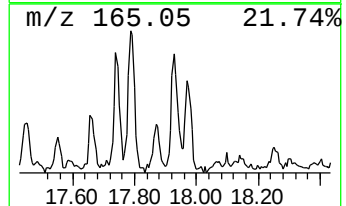
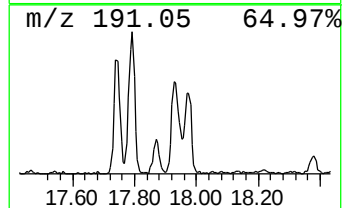
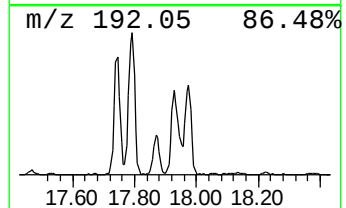
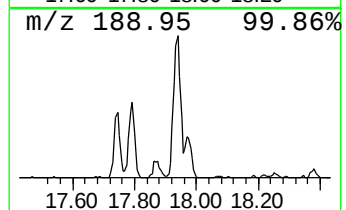
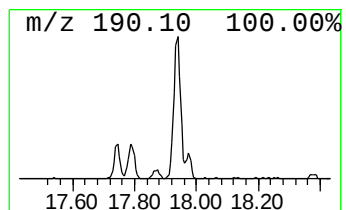
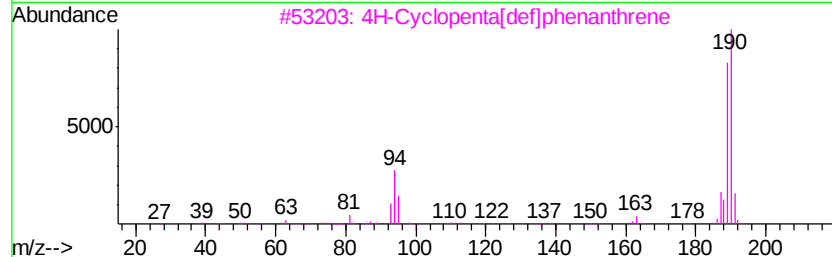
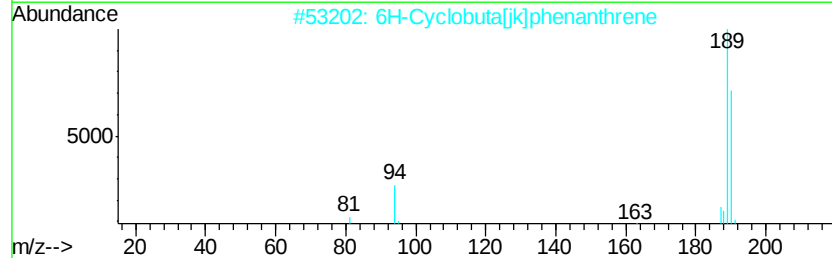
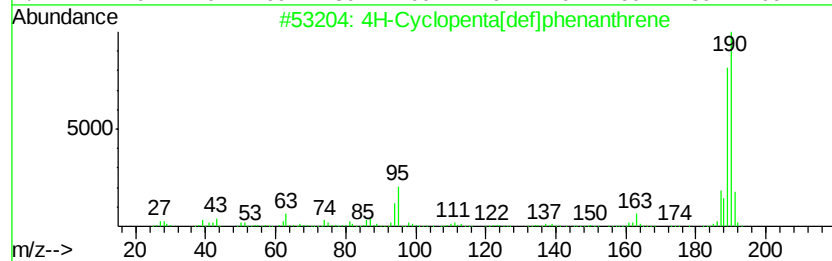
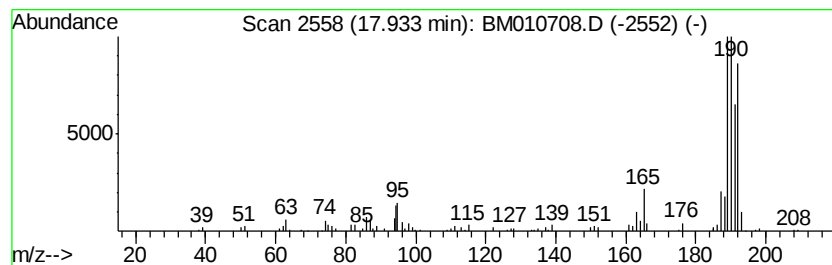
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Peak Number 3 4H-Cyclopenta[def]phenanthrene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.	
17.93	4.82 ng/ul	408515	Phenanthrene-d10	16.87	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	55
2	6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	46
3	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	46
4	5H-Dibenzo[a,d]cycloheptene	192	C15H12	000256-81-5	46
5	8,9-Dihydrocyclopenta[def]phenan...	192	C15H12	027410-55-5	42



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM062317\
Data File : BM010708.D
Acq On : 24 Jun 2017 02:02
Operator : SJ/JU
Sample : I3625-18
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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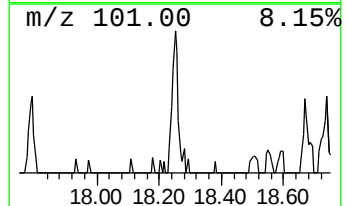
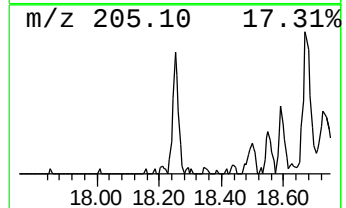
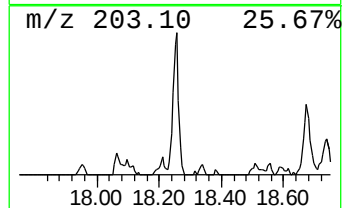
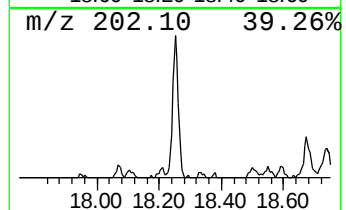
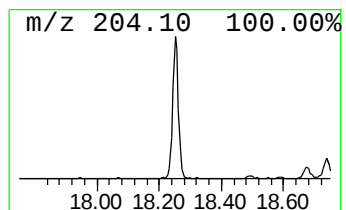
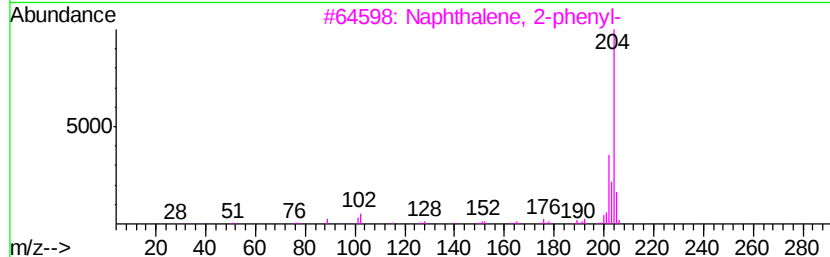
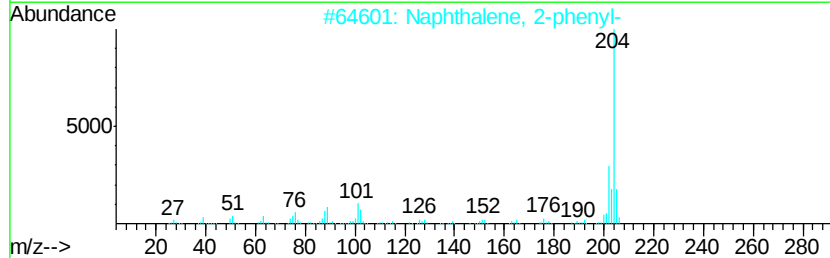
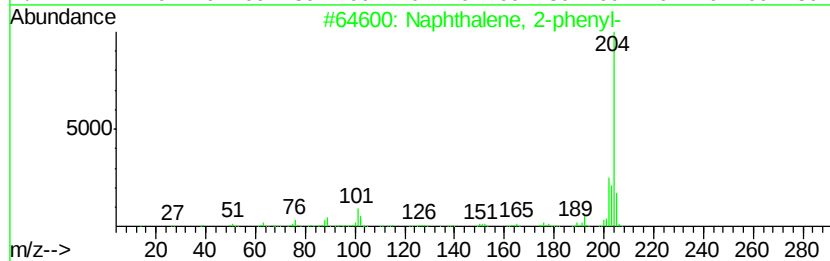
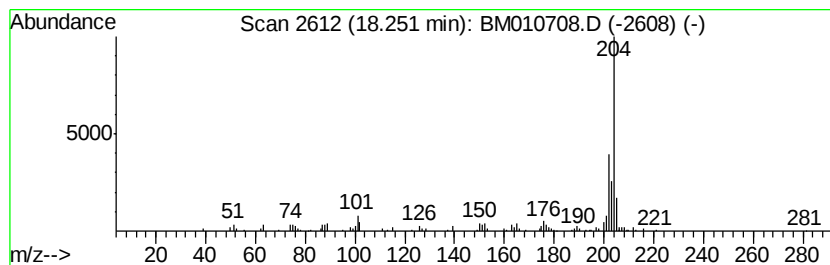
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Naphthalene, 2-phenyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.25	2.36 ng/ul	200510	Phenanthrene-d10	16.87

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	86
2		Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	83
3		Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	70
4		Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	70
5		Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	62



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM062317\
Data File : BM010708.D
Acq On : 24 Jun 2017 02:02
Operator : SJ/JU
Sample : I3625-18
Misc :
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Instrument :
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ClientSampleId :
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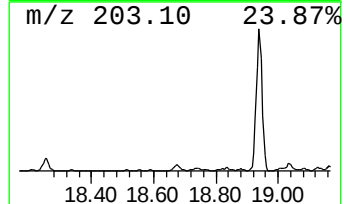
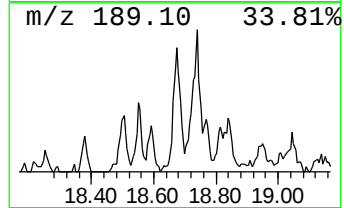
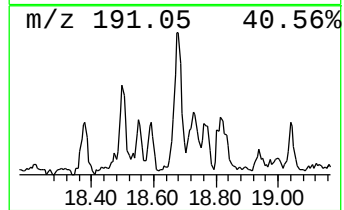
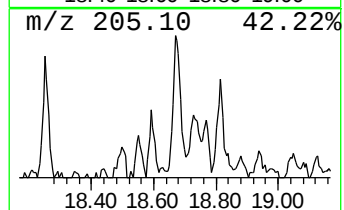
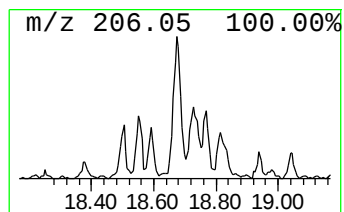
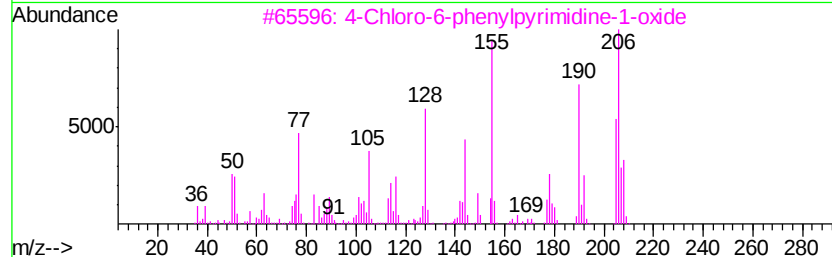
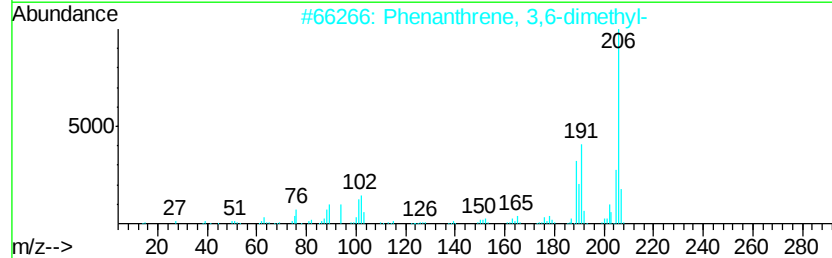
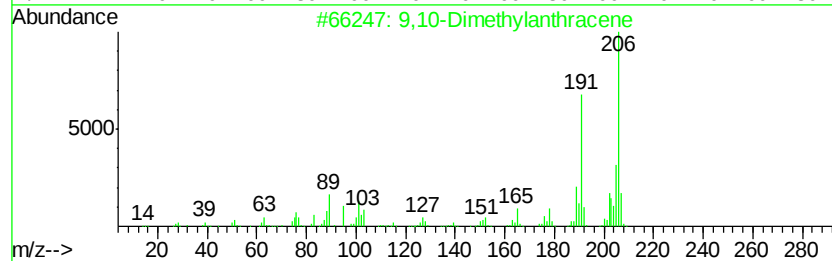
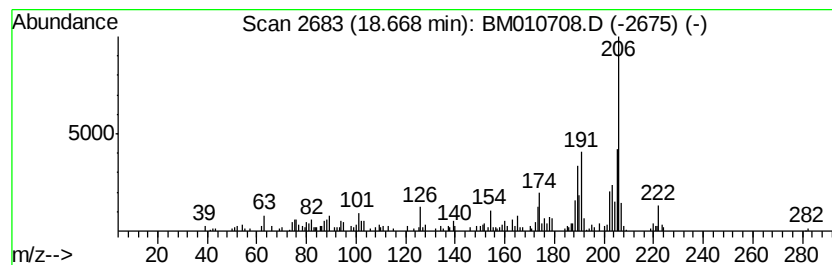
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 9,10-Dimethylantracene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.67	3.51 ng/ul	297516	Phenanthrene-d10	16.87

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	9,10-Dimethylantracene			206	C16H14	000781-43-1	92
2	Phenanthrene, 3,6-dimethyl-			206	C16H14	001576-67-6	89
3	4-Chloro-6-phenylpyrimidine-1-oxide			206	C10H7ClN2O	052816-77-0	83
4	Naphthalene, 1,2-dihydro-4-phenyl-			206	C16H14	007469-40-1	78
5	Phenanthrene, 2,7-dimethyl-			206	C16H14	001576-69-8	76



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM062317\
Data File : BM010708.D
Acq On : 24 Jun 2017 02:02
Operator : SJ/JU
Sample : I3625-18
Misc :
ALS Vial : 25 Sample Multiplier: 1

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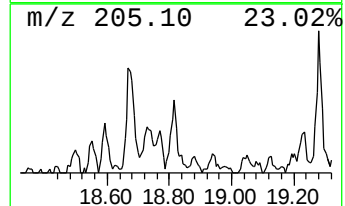
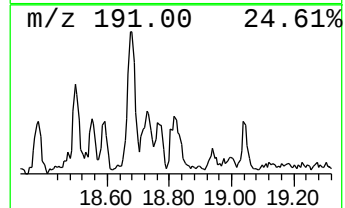
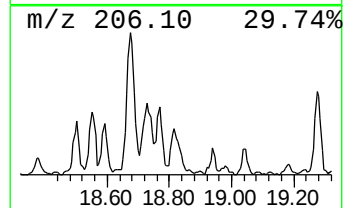
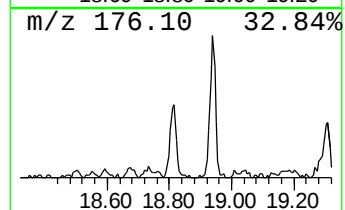
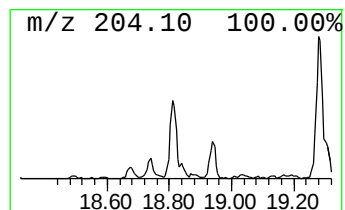
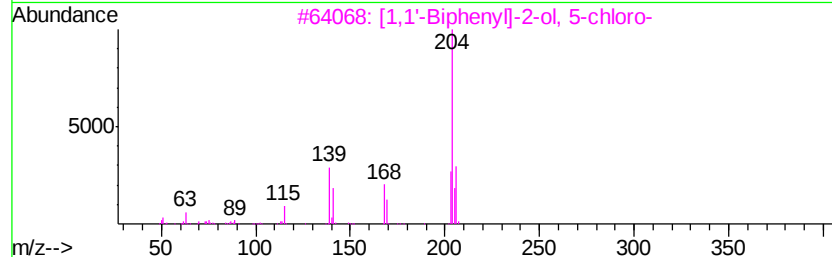
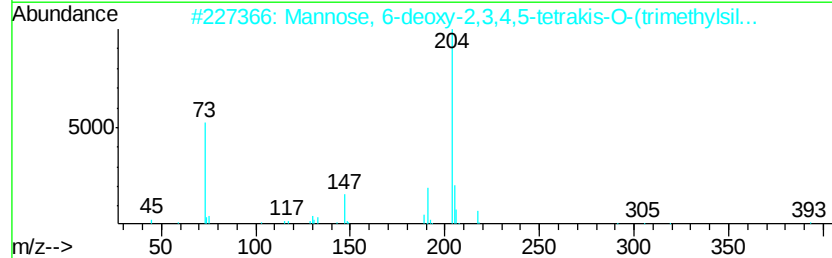
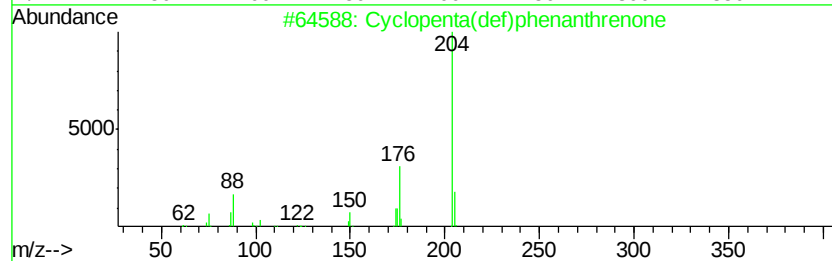
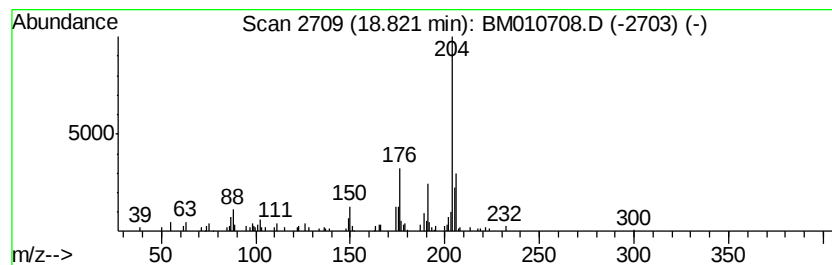
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Cyclopenta(def)phenanthrenone Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.82	2.20 ng/ul	186840	Phenanthrene-d10	16.87

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopenta(def)phenanthrenone	204	C15H8O	005737-13-3	68
2		Mannose, 6-deoxy-2,3,4,5-tetraki...	452	C18H44O5Si4	019127-15-2	47
3		[1,1'-Biphenyl]-2-ol, 5-chloro-	204	C12H9ClO	000607-12-5	43
4		Quinoline, 8-methoxy-5-nitro-	204	C10H8N2O3	1000298-88-7	43
5		[1,1'-Biphenyl]-4,4'-dicarbonitrile	204	C14H8N2	001591-30-6	38



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM062317\
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Acq On : 24 Jun 2017 02:02
Operator : SJ/JU
Sample : I3625-18
Misc :
ALS Vial : 25 Sample Multiplier: 1

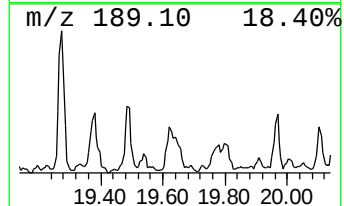
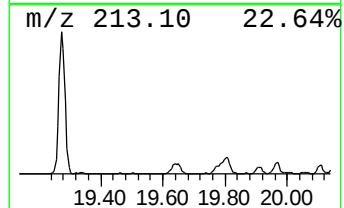
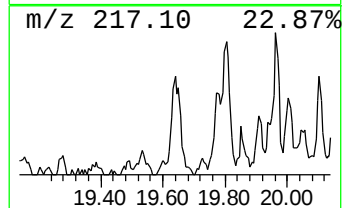
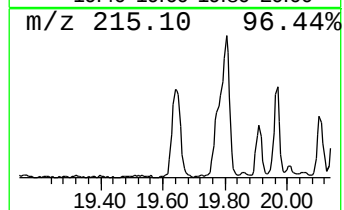
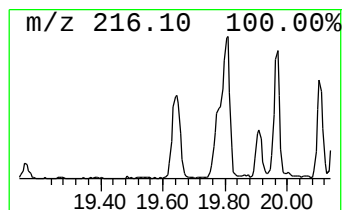
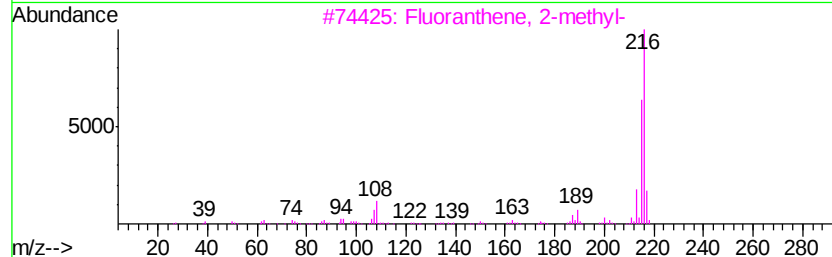
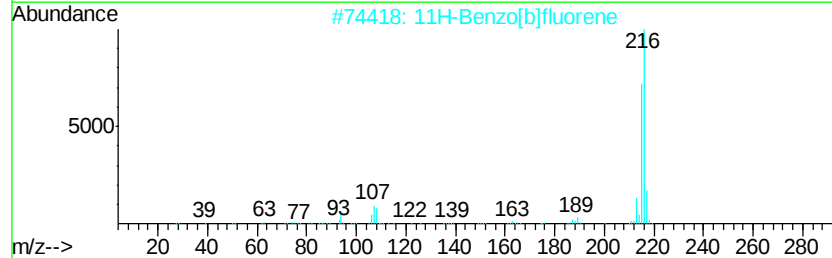
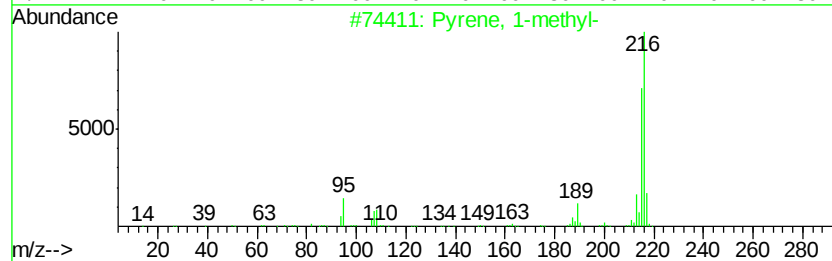
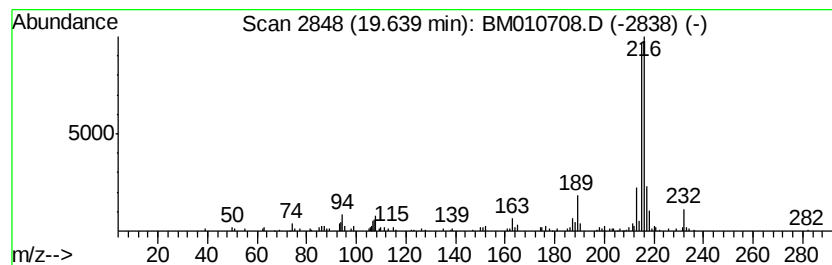
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Pyrene, 1-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.64	2.29 ng/ul	401001	Chrysene-d12	21.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Pyrene, 1-methyl-	216 C17H12	002381-21-7 90
2		11H-Benzo[b]fluorene	216 C17H12	000243-17-4 87
3		Fluoranthene, 2-methyl-	216 C17H12	033543-31-6 81
4		11H-Benzo[b]fluorene	216 C17H12	000243-17-4 76
5		11H-Benzo[a]fluorene	216 C17H12	000238-84-6 72



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM062317\
Data File : BM010708.D
Acq On : 24 Jun 2017 02:02
Operator : SJ/JU
Sample : I3625-18
Misc :
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Instrument :
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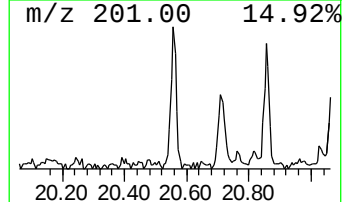
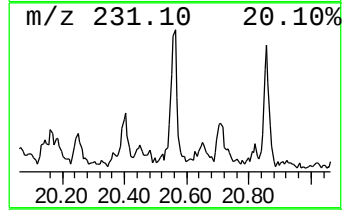
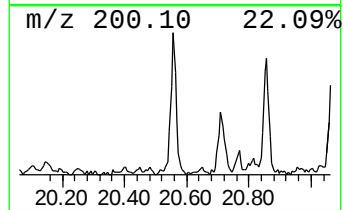
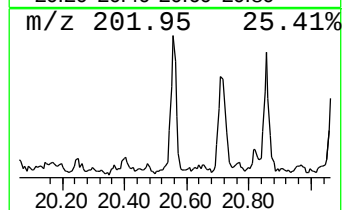
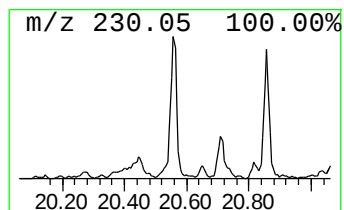
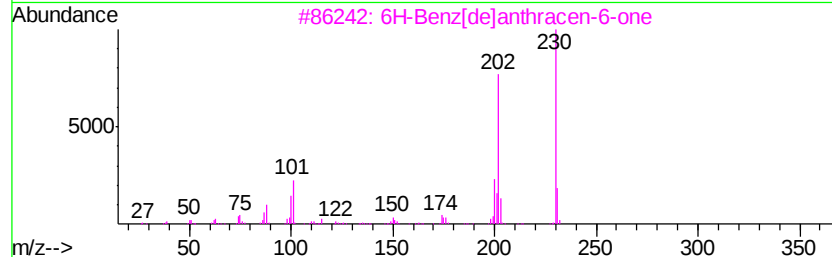
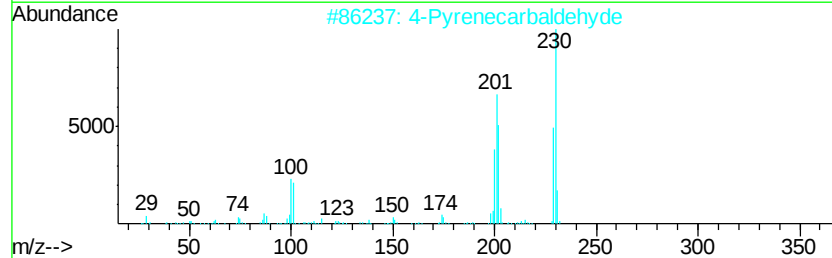
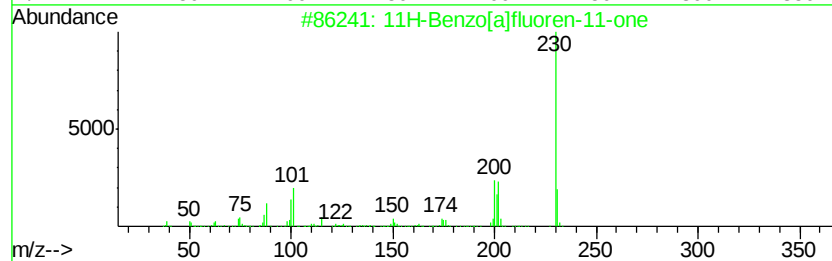
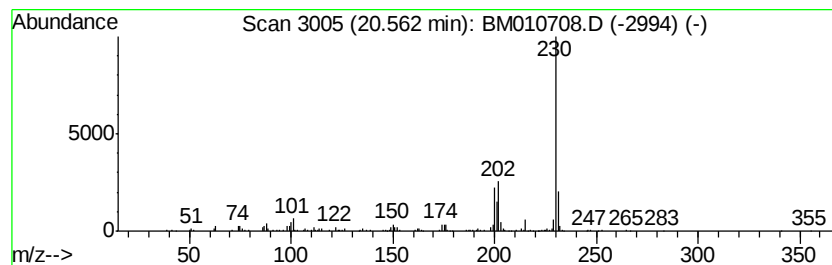
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 11H-Benzo[a]fluoren-11-one Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.56	2.41 ng/ul	422196	Chrysene-d12	21.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	11H-Benzo[a]fluoren-11-one	230	C17H10O	000479-79-8	96
2		4-Pyrenecarbaldehyde	230	C17H10O	022245-51-8	90
3		6H-Benz[de]anthracen-6-one	230	C17H10O	080252-14-8	89
4		7H-Benz[de]anthracen-7-one	230	C17H10O	000082-05-3	87
5		1-Pyrenecarboxaldehyde	230	C17H10O	003029-19-4	83



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM062317\
Data File : BM010708.D
Acq On : 24 Jun 2017 02:02
Operator : SJ/JU
Sample : I3625-18
Misc :
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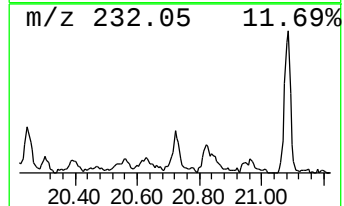
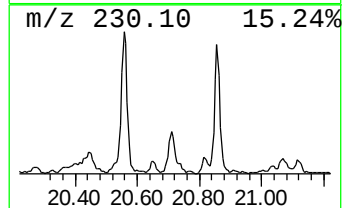
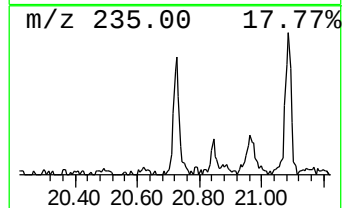
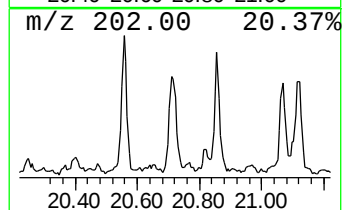
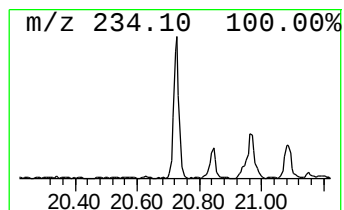
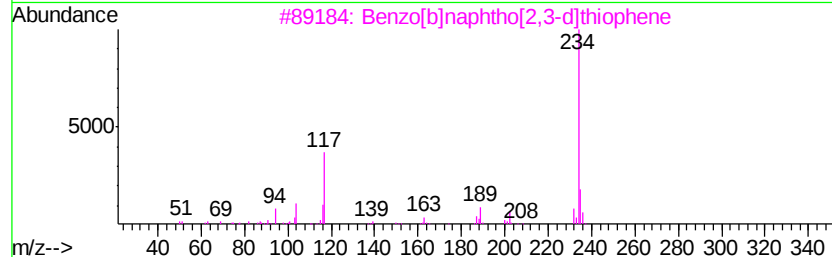
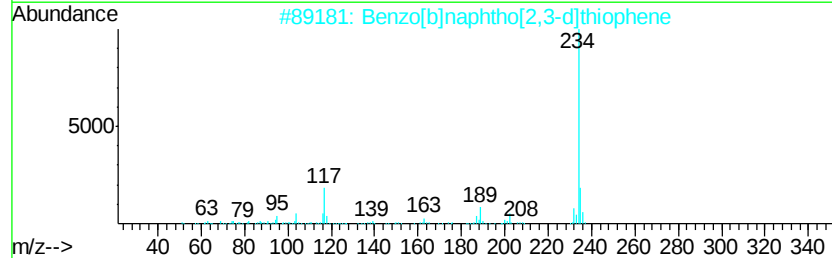
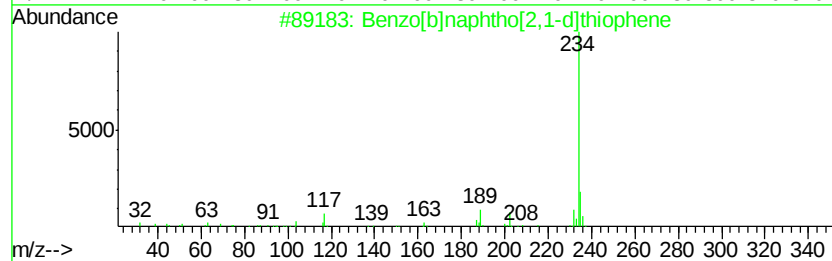
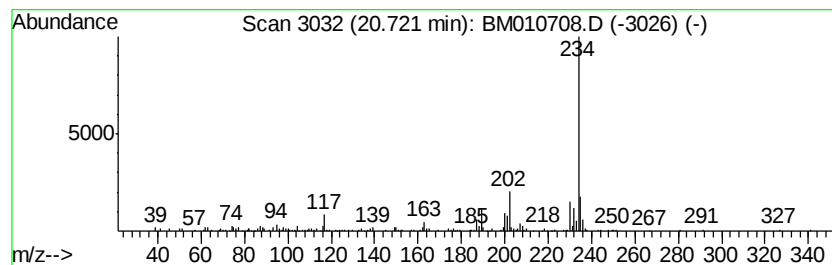
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 Benzo[b]naphtho[2,1-d]thiop... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.72	2.12 ng/ul	370325	Chrysene-d12	21.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzo[b]naphtho[2,1-d]thiophene	234	C16H10S	000239-35-0	95
2		Benzo[b]naphtho[2,3-d]thiophene	234	C16H10S	000243-46-9	89
3		Benzo[b]naphtho[2,3-d]thiophene	234	C16H10S	000243-46-9	76
4		Benzo[b]naphtho[2,1-d]thiophene	234	C16H10S	000239-35-0	76
5		Anthra(1,2-b)thiophene	234	C16H10S	000227-86-1	70



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM062317\
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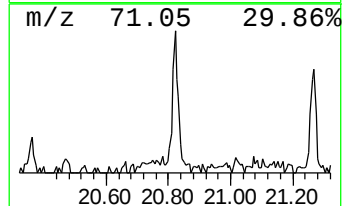
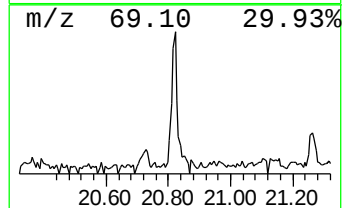
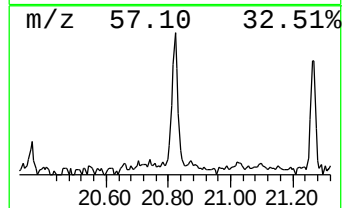
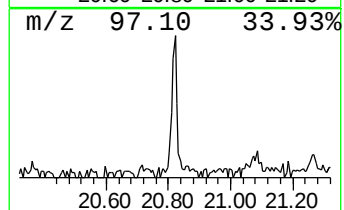
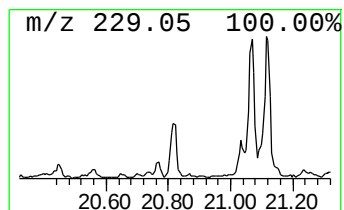
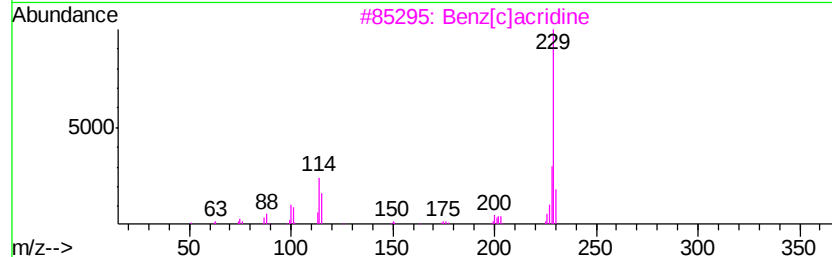
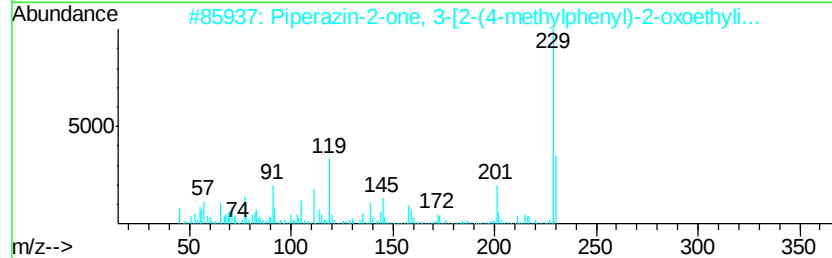
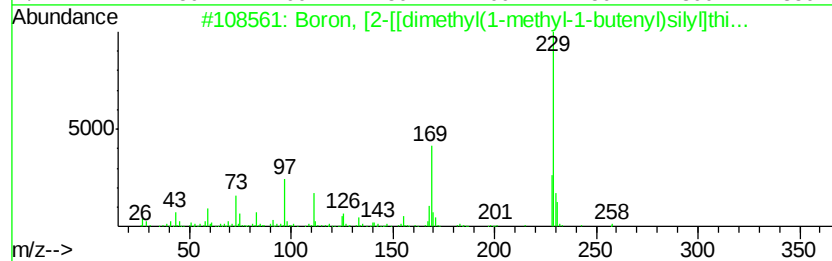
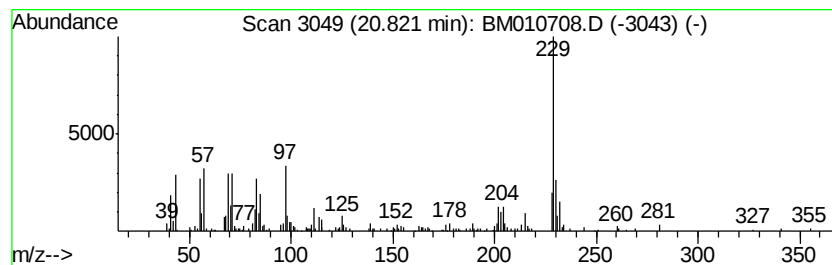
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 Boron, [2-[[dimethyl(1-meth... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.		
20.82	2.53 ng/ul	442559	Chrysene-d12	21.08		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Boron, [2-[[dimethyl(1-methyl-1-...	258	C11H23BS2Si	127855-37-2	58	
2	Piperazin-2-one, 3-[2-(4-methylp...	230	C13H14N2O2	063656-21-3	49	
3	Benz[c]acridine	229	C17H11N	000225-51-4	47	
4	Thiazole, 2-amino-4-(1H-indol-3-...	229	C12H11N3S	1000304-18-3	47	
5	Malononitrile, 2-indeno[1,2-b]py...	229	C15H7N3	1000316-60-9	43	



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM062317\
Data File : BM010708.D
Acq On : 24 Jun 2017 02:02
Operator : SJ/JU
Sample : I3625-18
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
F5A00

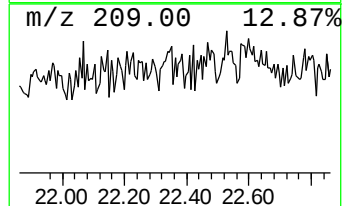
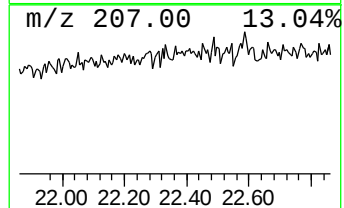
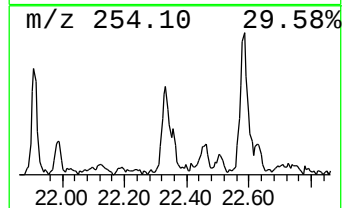
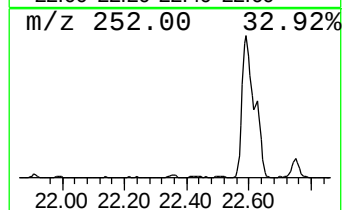
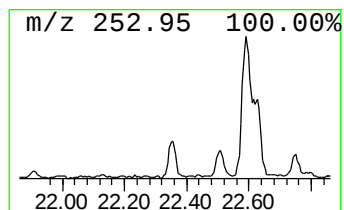
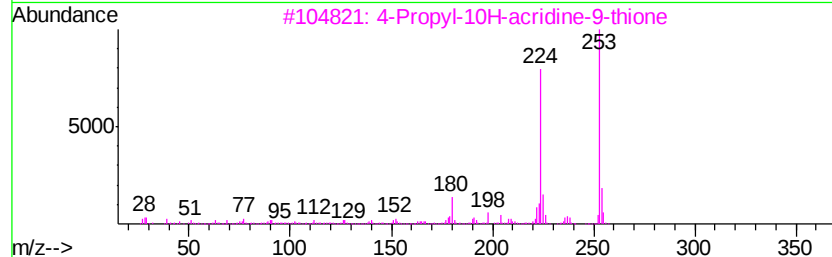
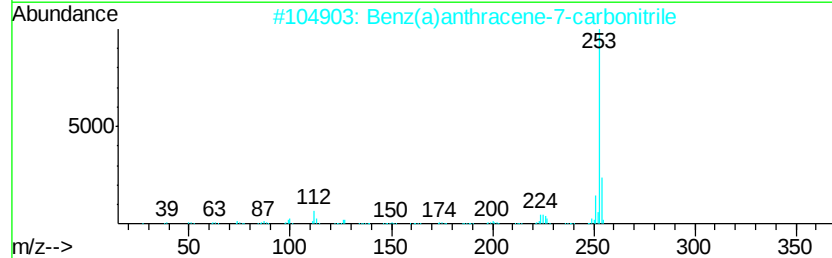
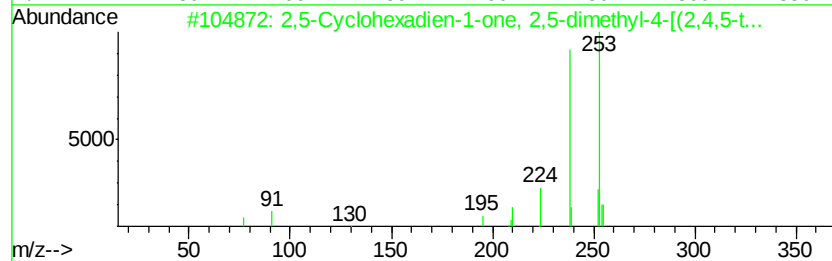
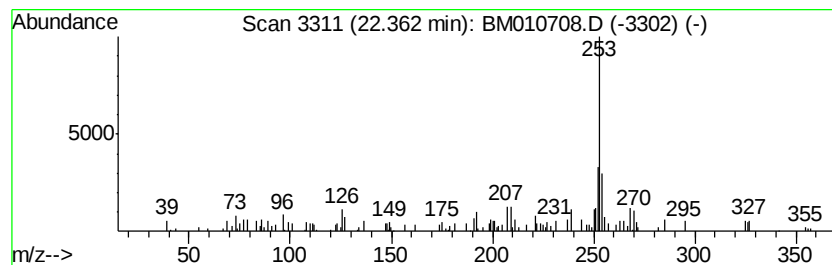
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 2,5-Cyclohexadien-1-one, 2,... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.36	2.36 ng/ul	150443	Perylene-d12	23.22

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2,5-Cyclohexadien-1-one, 2,5-dim...	253	C17H19NO	054245-93-1	53
2		Benz(a)anthracene-7-carbonitrile	253	C19H11N	007476-08-6	52
3		4-Propyl-10H-acridine-9-thione	253	C16H15NS	1000211-07-8	47
4		Acridin-1(2H)-one, 3,4-dihydro-3...	282	C18H22N2O	146352-02-5	43
5		6H-Indolo[2,3-b]quinoxaline, 9-c...	253	C14H8ClN3	1000304-55-7	43



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM062317\
Data File : BM010708.D
Acq On : 24 Jun 2017 02:02
Operator : SJ/JU
Sample : I3625-18
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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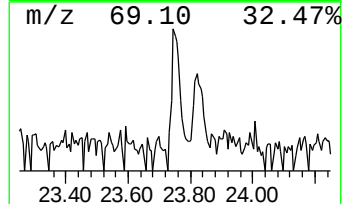
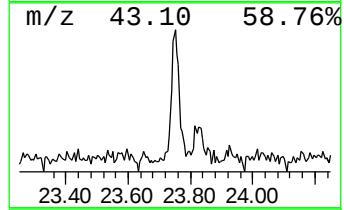
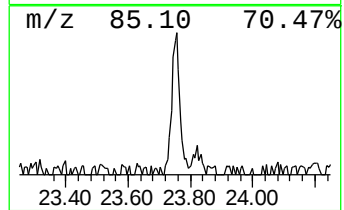
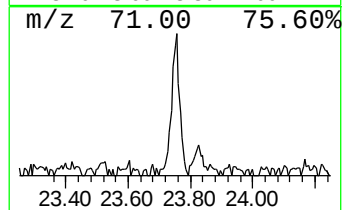
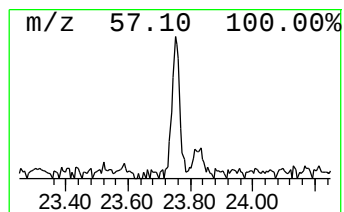
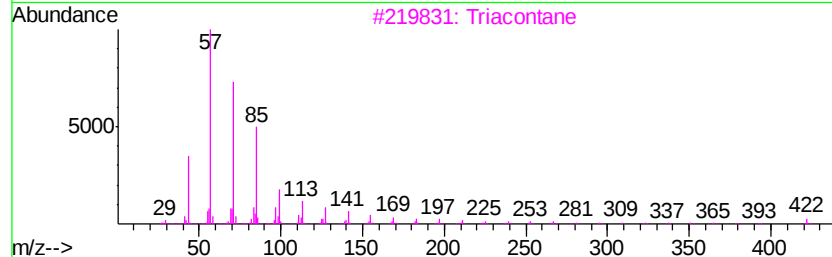
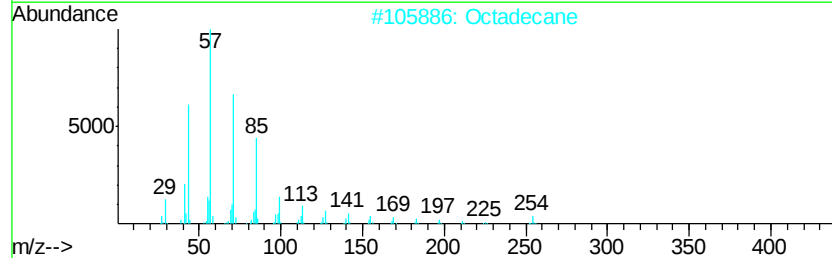
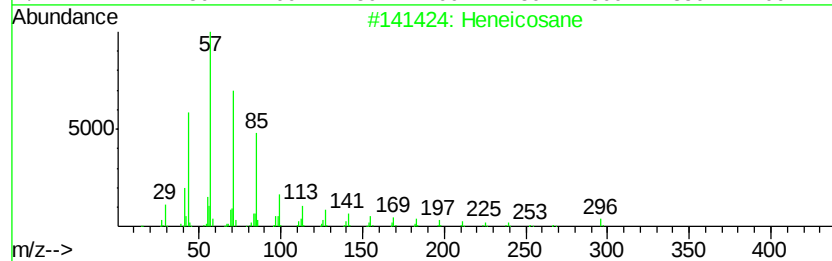
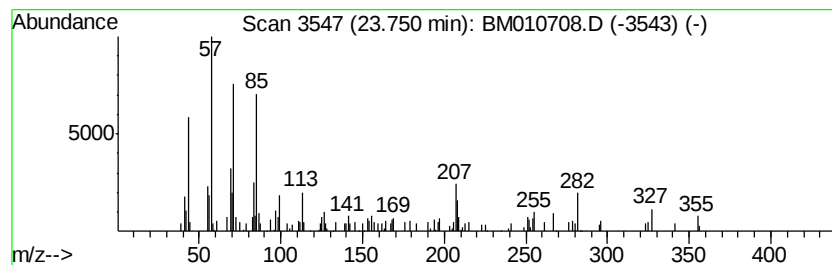
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Quant Title : SVOA CALIBRATION

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

Peak Number 14 (DEL) Alkane: Straight-Chai... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.75	2.04 ng/ul	129835	Perylene-d12	23.22

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heneicosane	296	C21H44	000629-94-7	78
2			Octadecane	254	C18H38	000593-45-3	60
3			Triacotane	422	C30H62	000638-68-6	53
4			Heneicosane	296	C21H44	000629-94-7	53
5			Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	49



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM062317\
 Data File : BM010708.D
 Acq On : 24 Jun 2017 02:02
 Operator : SJ/JU
 Sample : I3625-18
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 F5A00

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Phenanthrene, 1-m...	17.74	3.1	ng/ul	262482	4	16.87	1695800	20.0
1H-Cyclopropa[1]p...	17.79	6.4	ng/ul	545767	4	16.87	1695800	20.0
4H-Cyclopenta[def...	17.93	4.8	ng/ul	408515	4	16.87	1695800	20.0
Naphthalene, 2-ph...	18.25	2.4	ng/ul	200510	4	16.87	1695800	20.0
9,10-Dimethylanthe...	18.67	3.5	ng/ul	297516	4	16.87	1695800	20.0
Cyclopenta(def)ph...	18.82	2.2	ng/ul	186840	4	16.87	1695800	20.0
Pyrene, 1-methyl-	19.64	2.3	ng/ul	401001	5	21.08	3499050	20.0
11H-Benzo[a]fluor...	20.56	2.4	ng/ul	422196	5	21.08	3499050	20.0
Benzo[b]naphtho[2...	20.72	2.1	ng/ul	370325	5	21.08	3499050	20.0
Boron, [2-[[dimet...	20.82	2.5	ng/ul	442559	5	21.08	3499050	20.0
2,5-Cyclohexadien...	22.36	2.4	ng/ul	150443	6	23.22	1275650	20.0
(DEL) Alkane: Str...	23.75	2.0	ng/ul	129835	6	23.22	1275650	20.0