

Data Path : Z:\HPCHEM1\BNA M\DATA\BM030518\
 Data File : BM014485.D
 Acq On : 05 Mar 2018 22:35
 Operator : SJ/JU
 Sample : J1678-09DL 5X
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BE5T9DL

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-BM021418.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	6.740	633	638	652	rVV5	50516	147662	5.24%	0.499%
2	6.869	655	660	668	rBV	61788	109234	3.88%	0.369%
3	7.063	686	693	702	rBV	87464	180793	6.42%	0.611%
4	7.510	762	769	784	rBV	418894	724276	25.70%	2.448%
5	8.269	891	898	908	rBV4	55381	155165	5.51%	0.524%
6	8.687	961	969	979	rBV2	79398	179058	6.35%	0.605%
7	9.398	1084	1090	1096	rBV4	60645	119818	4.25%	0.405%
8	9.975	1182	1188	1196	rBV7	47053	145135	5.15%	0.490%
9	10.281	1233	1240	1247	rBV	464098	892171	31.66%	3.015%
10	10.475	1268	1273	1274	rBV3	25275	31274	1.11%	0.106%
11	11.969	1524	1527	1536	rBV6	11222	28233	1.00%	0.095%
12	12.204	1559	1567	1571	rBV6	13817	28866	1.02%	0.098%
13	13.492	1781	1786	1791	rBV4	23305	42746	1.52%	0.144%
14	13.598	1799	1804	1809	rBV2	165611	275412	9.77%	0.931%
15	13.645	1809	1812	1819	rVV2	35084	63186	2.24%	0.214%
16	13.863	1843	1849	1862	rBV3	225046	481354	17.08%	1.627%
17	14.075	1880	1885	1891	rBV7	16711	30489	1.08%	0.103%
18	14.175	1896	1902	1909	rVV	701577	1128861	40.06%	3.815%
19	14.239	1909	1913	1923	rVB2	75880	135164	4.80%	0.457%
20	14.586	1967	1972	1978	rVV	68087	118513	4.21%	0.400%
21	14.663	1981	1985	1995	rVB7	20513	41296	1.47%	0.140%
22	14.798	2003	2008	2014	rBV5	16640	29634	1.05%	0.100%
23	14.975	2029	2038	2042	rBV5	18243	46006	1.63%	0.155%
24	15.033	2042	2048	2053	rVB6	20250	41470	1.47%	0.140%
25	15.180	2068	2073	2078	rBV	272291	415137	14.73%	1.403%
26	15.239	2078	2083	2093	rVB3	81008	167475	5.94%	0.566%
27	15.533	2129	2133	2141	rVB5	38042	69434	2.46%	0.235%
28	15.698	2153	2161	2166	rVB6	30520	67578	2.40%	0.228%
29	16.251	2247	2255	2260	rBV7	32579	66493	2.36%	0.225%
30	16.292	2260	2262	2267	rVB4	23306	32461	1.15%	0.110%
31	16.474	2289	2293	2296	rBV5	25913	28831	1.02%	0.097%
32	16.516	2296	2300	2304	rVB4	26139	33331	1.18%	0.113%
33	16.610	2310	2316	2321	rBV2	59951	91708	3.25%	0.310%
34	16.680	2323	2328	2333	rVV7	34513	76999	2.73%	0.260%

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35	16.757	2336	2341	2347	rVB	68284	100861	3.58%	0.341%
36	16.939	2366	2372	2375	rBV	820483	1208441	42.88%	4.084%
37	16.980	2375	2379	2385	rVV	1303999	1876609	66.59%	6.342%
38	17.039	2385	2389	2392	rVV2	313613	481655	17.09%	1.628%
39	17.074	2392	2395	2400	rVB	251405	349055	12.39%	1.180%
40	17.151	2405	2408	2411	rVB5	21042	29069	1.03%	0.098%
41	17.333	2435	2439	2440	rBV4	29049	34093	1.21%	0.115%
42	17.369	2440	2445	2450	rVB2	75804	148092	5.25%	0.500%
43	17.457	2457	2460	2463	rBV2	33146	39503	1.40%	0.133%
44	17.657	2490	2494	2495	rBV4	24980	29352	1.04%	0.099%
45	17.745	2505	2509	2513	rBV7	19899	31855	1.13%	0.108%
46	17.816	2515	2521	2525	rVV	199641	323016	11.46%	1.092%
47	17.863	2525	2529	2535	rVV	251466	395809	14.05%	1.338%
48	17.945	2539	2543	2548	rVV	84975	128882	4.57%	0.436%
49	18.010	2548	2554	2558	rVV3	242863	452229	16.05%	1.528%
50	18.045	2558	2560	2568	rVB	141571	197628	7.01%	0.668%
51	18.127	2570	2574	2580	rBV9	25682	49301	1.75%	0.167%
52	18.180	2580	2583	2586	rBV5	23492	32747	1.16%	0.111%
53	18.321	2603	2607	2613	rVB	127082	168507	5.98%	0.569%
54	18.386	2613	2618	2625	rBV	140836	210260	7.46%	0.711%
55	18.568	2642	2649	2655	rBV	64969	156670	5.56%	0.529%
56	18.621	2655	2658	2661	rVB4	44023	52181	1.85%	0.176%
57	18.663	2663	2665	2670	rVB6	43797	53508	1.90%	0.181%
58	18.745	2674	2679	2683	rBV	133340	222360	7.89%	0.751%
59	18.810	2688	2690	2693	rVB4	37550	36908	1.31%	0.125%
60	18.904	2699	2706	2711	rBV3	127830	271941	9.65%	0.919%
61	19.010	2719	2724	2731	rBV	2126675	2818141	100.00%	9.523%
62	19.080	2733	2736	2739	rBV5	31521	44025	1.56%	0.149%
63	19.115	2739	2742	2747	rBV7	48692	78005	2.77%	0.264%
64	19.257	2760	2766	2769	rVB3	80504	141329	5.01%	0.478%
65	19.351	2776	2782	2783	rBV2	641101	845167	29.99%	2.856%
66	19.380	2783	2787	2795	rVB	1785178	2440053	86.58%	8.246%
67	19.451	2796	2799	2803	rBV2	83929	112395	3.99%	0.380%
68	19.562	2815	2818	2822	rVB	85565	98942	3.51%	0.334%
69	19.633	2827	2830	2837	rVB	81133	125775	4.46%	0.425%
70	19.710	2837	2843	2850	rBV3	140879	364692	12.94%	1.232%
71	19.845	2863	2866	2867	rBV3	88230	96735	3.43%	0.327%

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72	19.880	2869	2872	2877	rVB2	185221	293521	10.42%	0.992%
73	19.980	2886	2889	2893	rBV	114695	183210	6.50%	0.619%
74	20.039	2895	2899	2903	rVB2	178078	277653	9.85%	0.938%
75	20.180	2920	2923	2927	rBV	109124	150932	5.36%	0.510%
76	20.227	2927	2931	2939	rVB2	136898	193690	6.87%	0.655%
77	20.439	2965	2967	2971	rVB5	72241	58159	2.06%	0.197%
78	20.645	2999	3002	3005	rVB	193692	210137	7.46%	0.710%
79	20.798	3025	3028	3031	rBV	167548	212789	7.55%	0.719%
80	20.833	3032	3034	3038	rVB	118675	119563	4.24%	0.404%
81	20.945	3050	3053	3058	rVB	180649	229434	8.14%	0.775%
82	21.151	3082	3088	3092	rBV2	1201730	2520843	89.45%	8.519%
83	21.192	3092	3095	3099	rVB	886333	990253	35.14%	3.346%
84	22.692	3345	3350	3355	rBV	620159	1321196	46.88%	4.465%
85	23.150	3424	3428	3432	rBV	351626	547375	19.42%	1.850%
86	23.245	3440	3444	3449	rVB2	566332	949247	33.68%	3.208%
87	23.339	3455	3460	3464	rBV	516584	862774	30.62%	2.916%

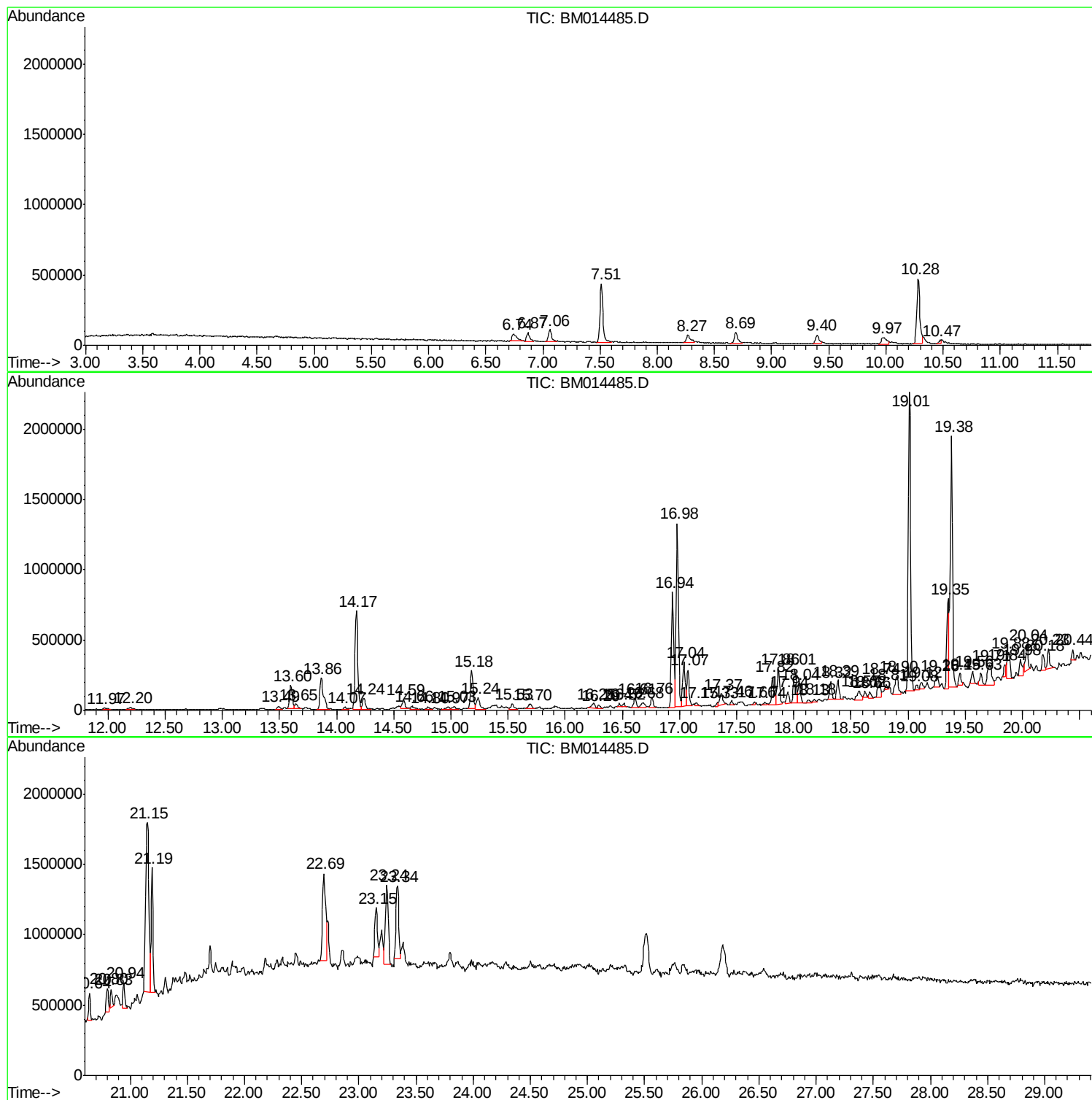
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Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM021418.M
Quant Title : SVOA CALIBRATION

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



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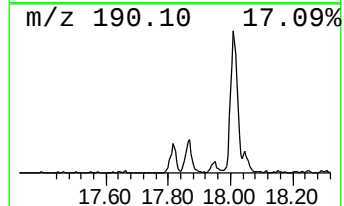
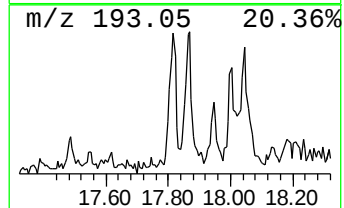
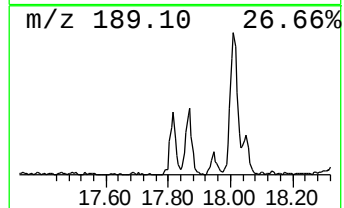
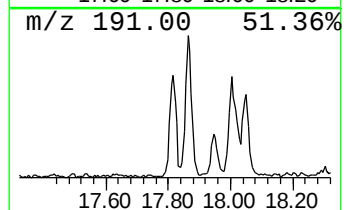
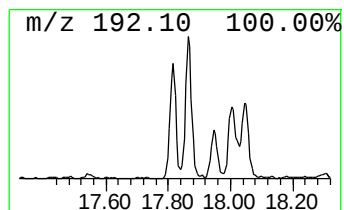
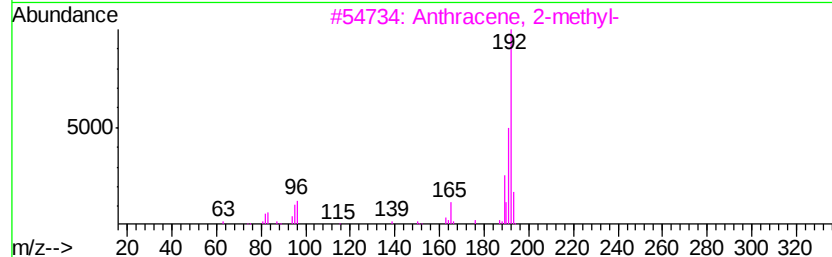
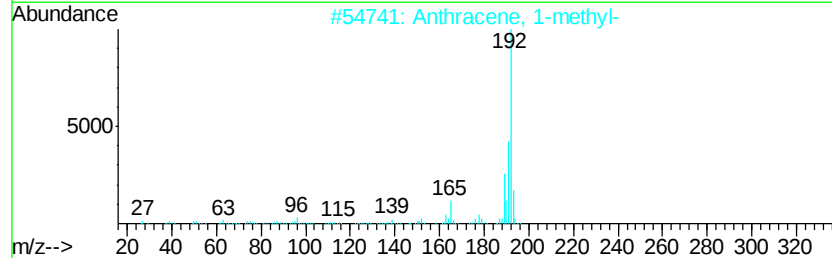
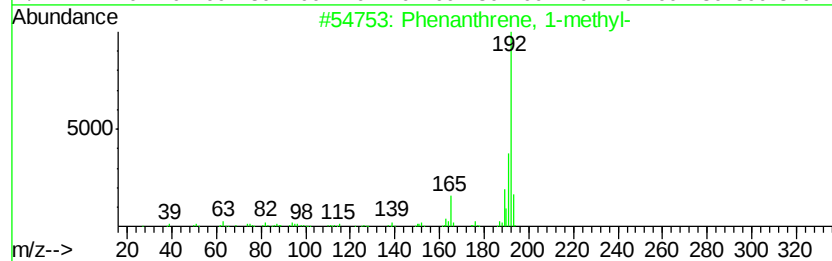
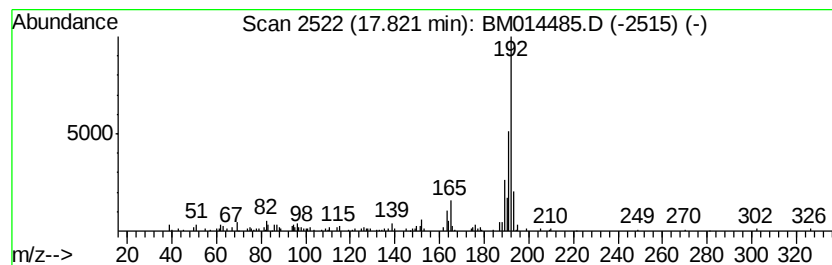
Instrument :
BNA_M
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Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM021418.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.	
17.82	5.35 ng/ul	323016	Phenanthrene-d10	16.94	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	95
2	Anthracene, 1-methyl-	192	C15H12	000610-48-0	95
3	Anthracene, 2-methyl-	192	C15H12	000613-12-7	91
4	1H-Cyclopropa[1]phenanthrene,1a,...	192	C15H12	000949-41-7	91
5	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	91



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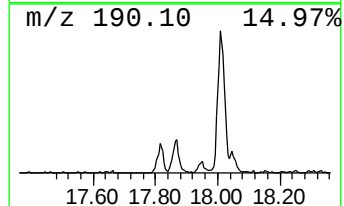
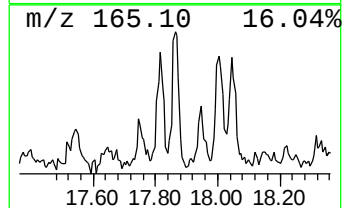
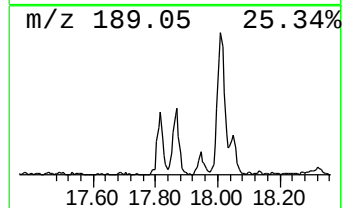
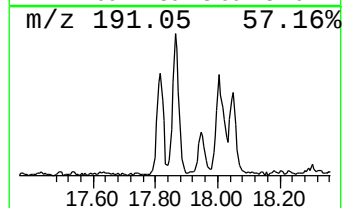
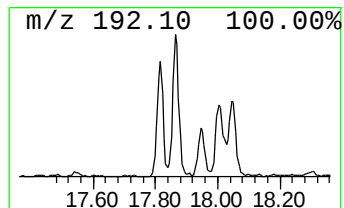
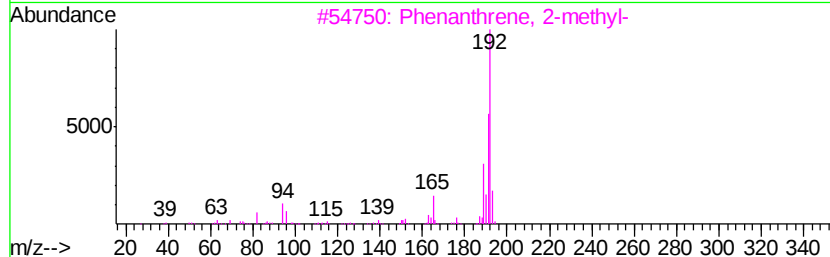
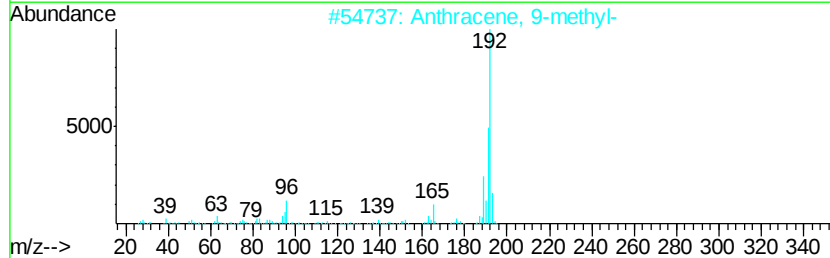
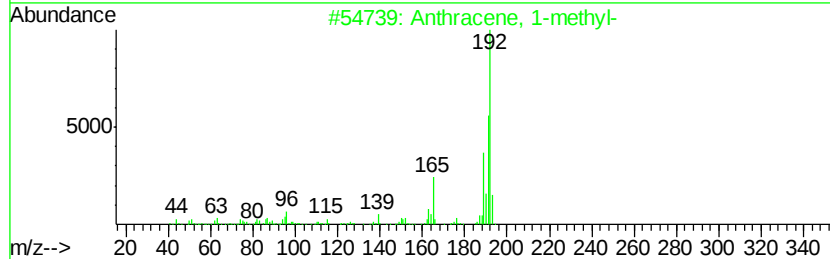
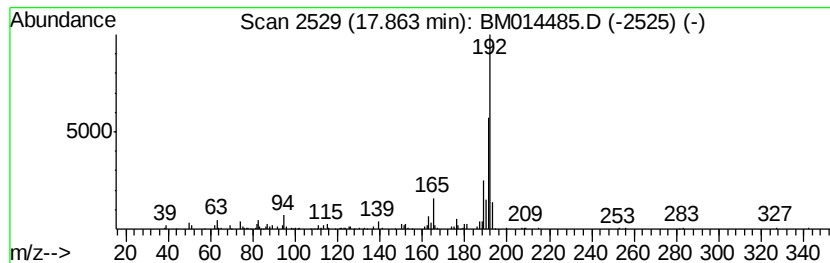
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.86	6.55 ng/ul	395809	Phenanthrene-d10	16.94

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Anthracene, 1-methyl-	192	C15H12	000610-48-0	93
2		Anthracene, 9-methyl-	192	C15H12	000779-02-2	93
3		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	93
4		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	93
5		Anthracene, 1-methyl-	192	C15H12	000610-48-0	90



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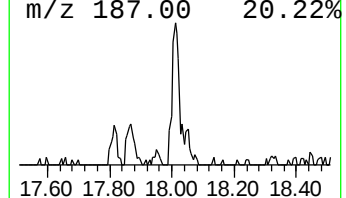
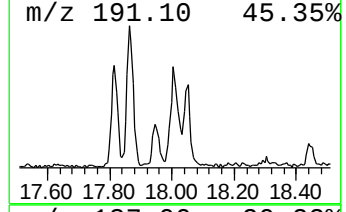
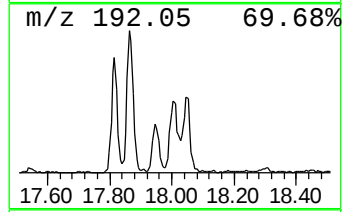
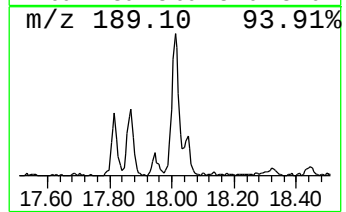
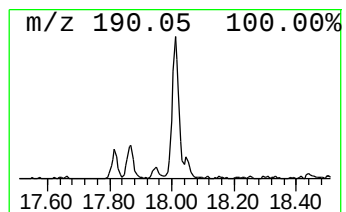
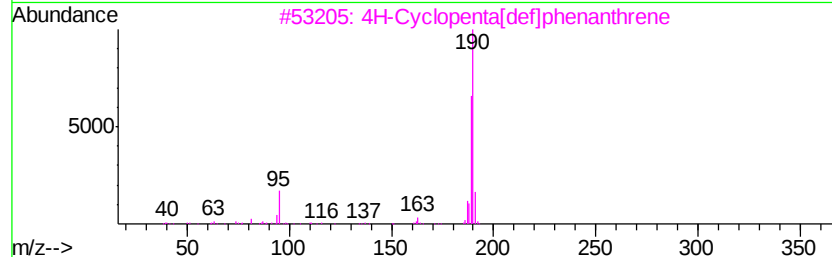
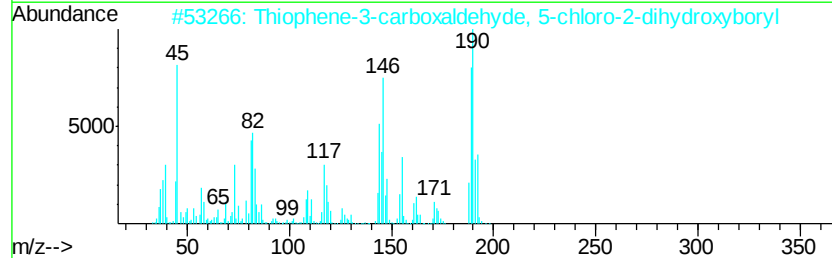
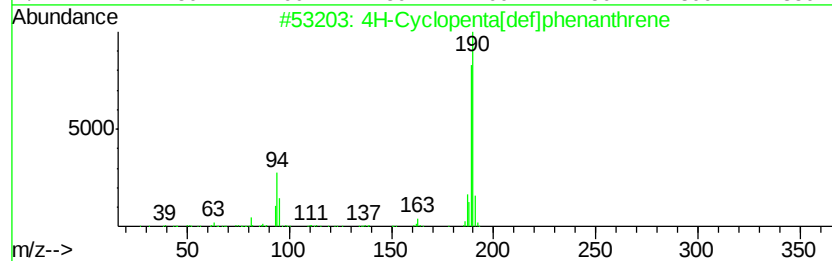
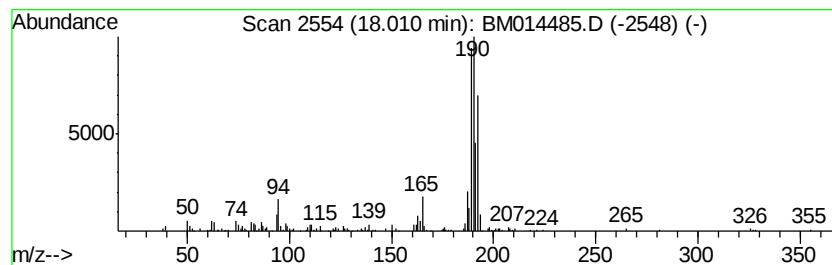
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 4H-Cyclopenta[def]phenanthrene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.01	7.48 ng/ul	452229	Phenanthrene-d10	16.94

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	62
2		Thiophene-3-carboxaldehyde, 5-ch...	190	C5H4BClO3S	036155-87-0	53
3		4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	52
4		4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	49
5		Carbonic acid, ethyl 2-formyl-4,...	262	C10H8Cl2O4	1000331-34-4	40



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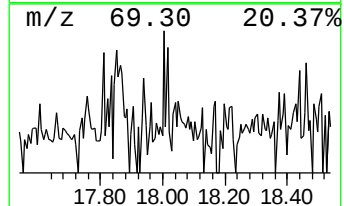
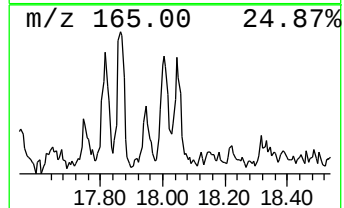
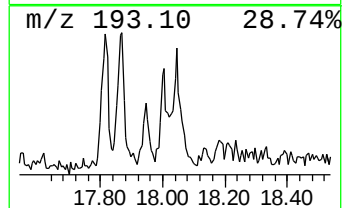
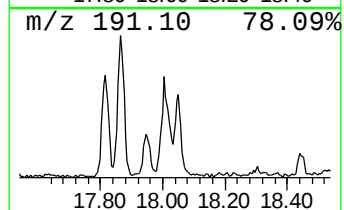
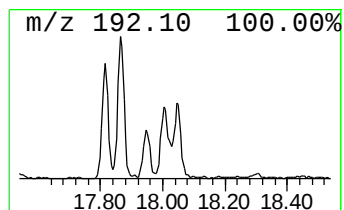
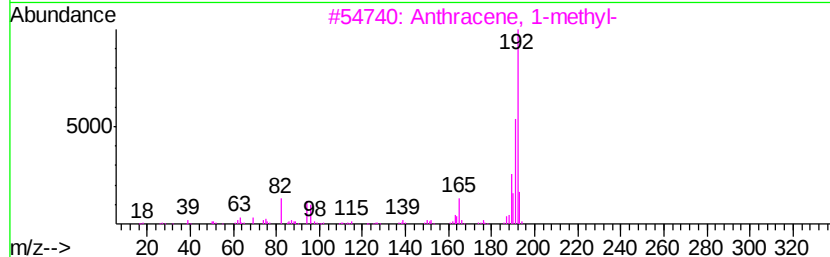
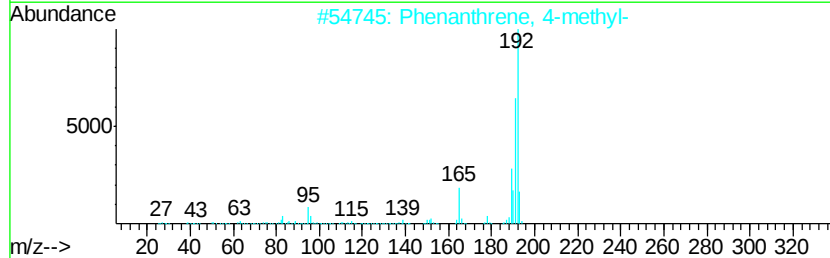
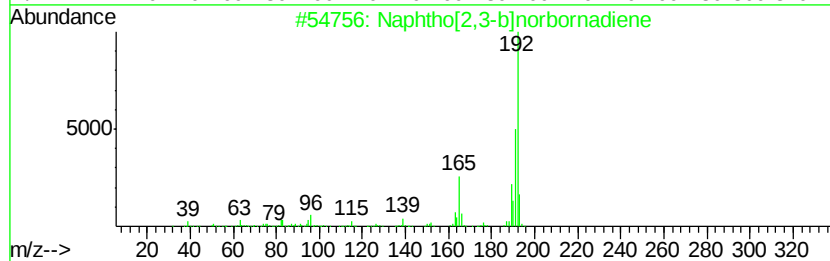
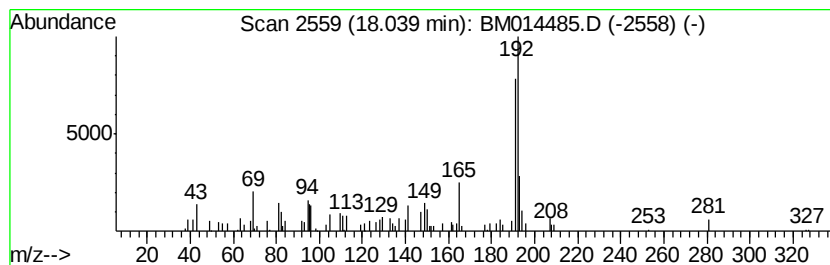
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Naphtho[2,3-b]norbornadiene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.04	3.27 ng/ul	197628	Phenanthrene-d10	16.94

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Naphtho[2,3-b]norbornadiene	192	C15H12	107426-38-0	64
2		Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	59
3		Anthracene, 1-methyl-	192	C15H12	000610-48-0	59
4		2-Hydroxy-4-methyl-6-(2-thienyl)...	192	C9H8N2OS	026963-45-1	58
5		Anthracene, 9-methyl-	192	C15H12	000779-02-2	58



Data Path : Z:\HPCHEM1\BNA M\DATA\BM030518\
Data File : BM014485.D
Acq On : 05 Mar 2018 22:35
Operator : SJ/JU
Sample : J1678-09DL 5X
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
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ClientSampleId :
BE5T9DL

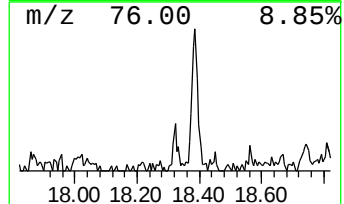
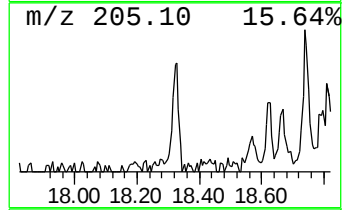
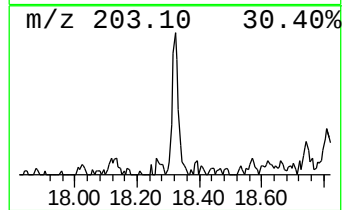
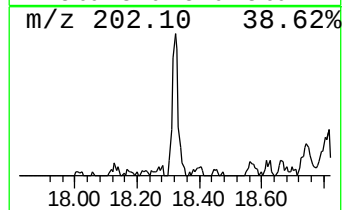
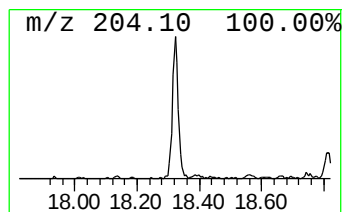
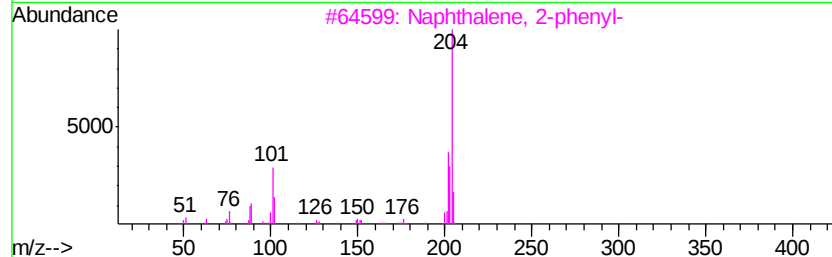
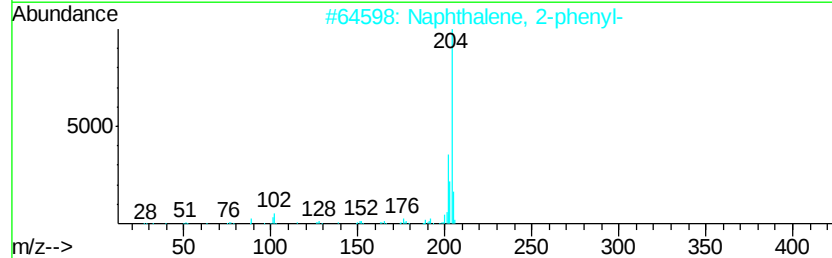
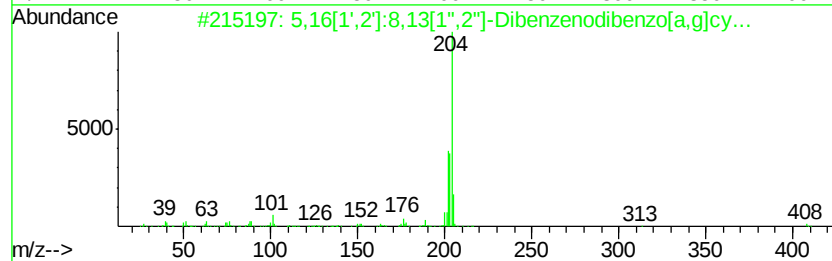
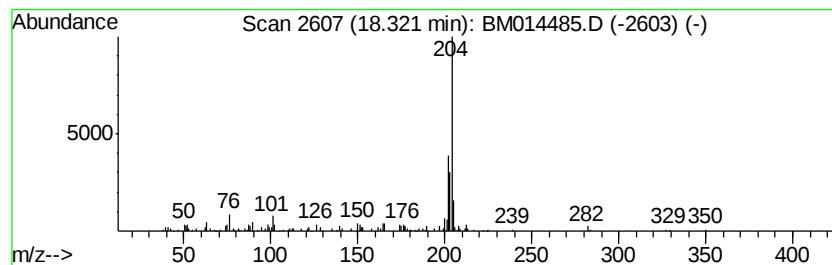
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 5,16[1',2']:8,13[1'',2'']-D... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.32	2.79 ng/ul	168507	Phenanthrene-d10	16.94

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	70
3	Naphthalene, 2-phenyl-	204	C16H12		000612-94-2	70
4	Tricyclo[8.2.2.2(4,7)]hexadeca-2...	204	C16H12		006572-60-7	62
5	Naphthalene, 2-phenyl-	204	C16H12		000612-94-2	58



Data Path : Z:\HPCHEM1\BNA M\DATA\BM030518\
Data File : BM014485.D
Acq On : 05 Mar 2018 22:35
Operator : SJ/JU
Sample : J1678-09DL 5X
Misc :
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ClientSampleId :
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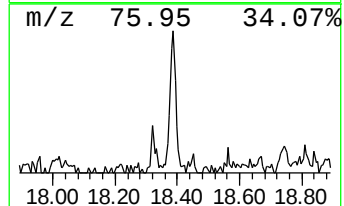
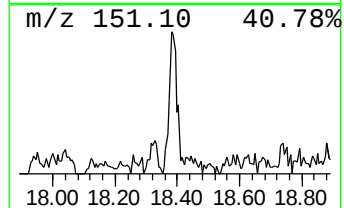
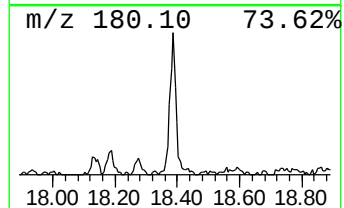
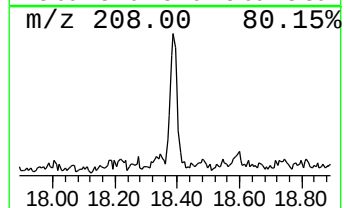
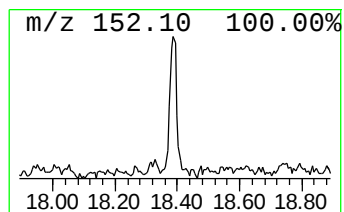
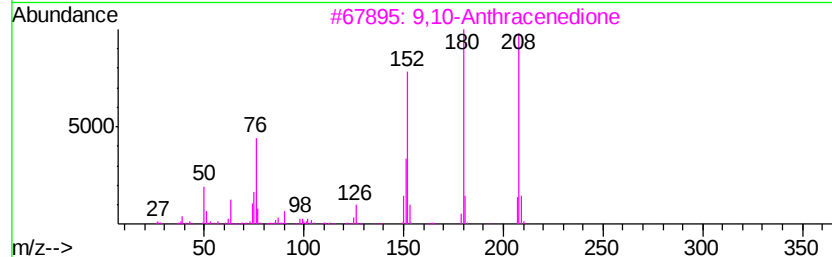
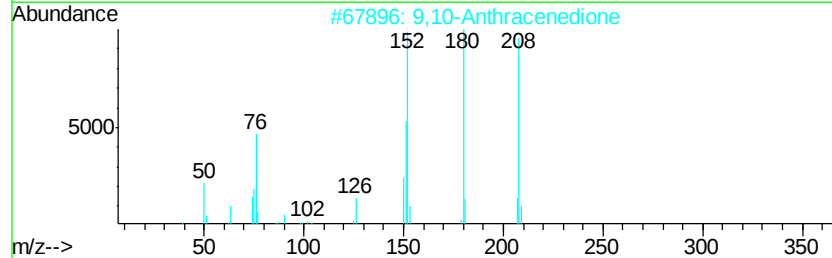
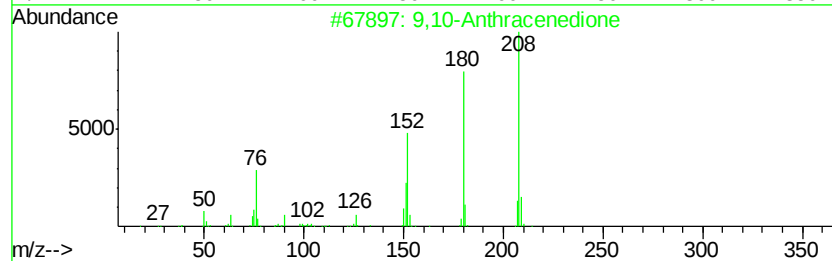
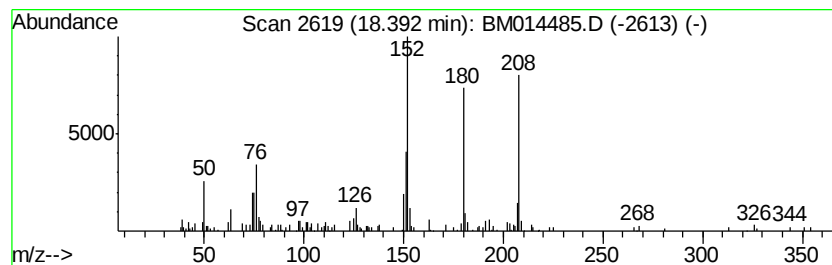
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 9,10-Anthracenedione Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.39	3.48 ng/ul	210260	Phenanthrene-d10	16.94

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	9,10-Anthracenedione		208	C14H8O2	000084-65-1	94
2	9,10-Anthracenedione		208	C14H8O2	000084-65-1	93
3	9,10-Anthracenedione		208	C14H8O2	000084-65-1	91
4	Benzo[ghi]cinnoline		180	C12H8N2	000230-31-9	70
5	9,10-Phenanthrenedione		208	C14H8O2	000084-11-7	70



Data Path : Z:\HPCHEM1\BNA M\DATA\BM030518\
Data File : BM014485.D
Acq On : 05 Mar 2018 22:35
Operator : SJ/JU
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ALS Vial : 18 Sample Multiplier: 1

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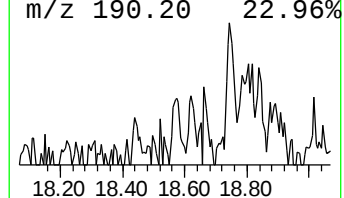
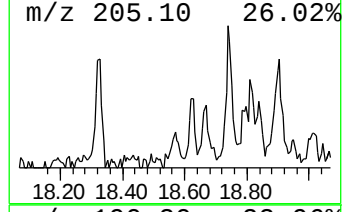
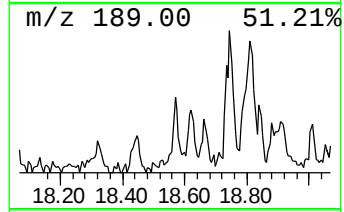
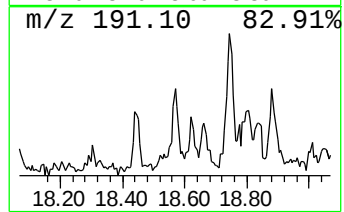
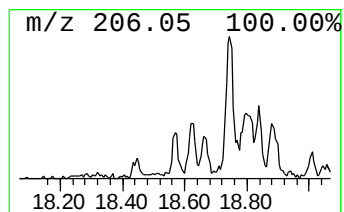
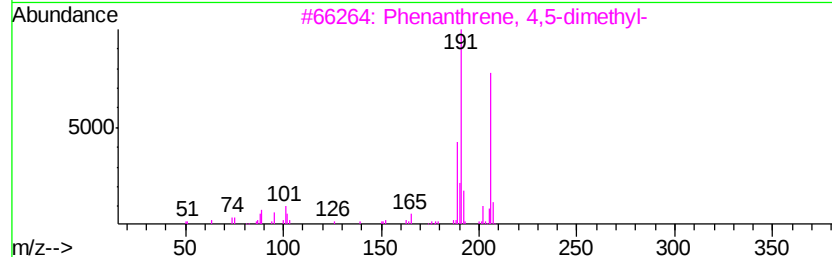
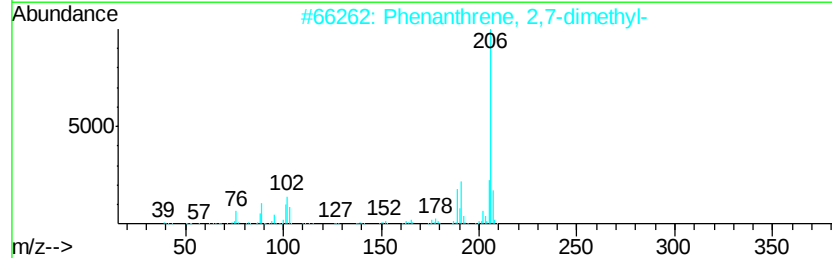
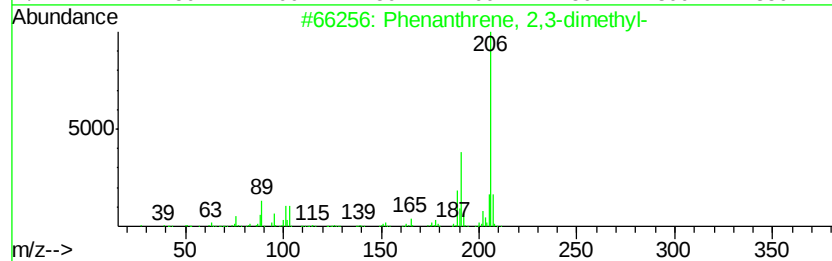
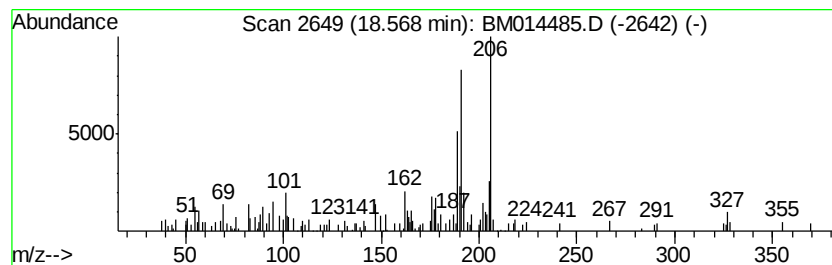
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Phenanthrene, 2,3-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.57	2.59 ng/ul	156670	Phenanthrene-d10	16.94

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phenanthrene, 2,3-dimethyl-	206	C16H14	003674-65-5	91
2		Phenanthrene, 2,7-dimethyl-	206	C16H14	001576-69-8	70
3		Phenanthrene, 4,5-dimethyl-	206	C16H14	003674-69-9	70
4		9,10-Dimethylantracene	206	C16H14	000781-43-1	68
5		9,10-Dimethylantracene	206	C16H14	000781-43-1	62



Data Path : Z:\HPCHEM1\BNA M\DATA\BM030518\
Data File : BM014485.D
Acq On : 05 Mar 2018 22:35
Operator : SJ/JU
Sample : J1678-09DL 5X
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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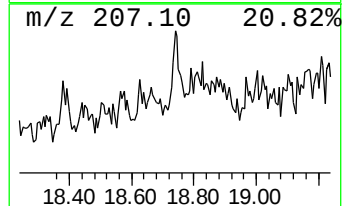
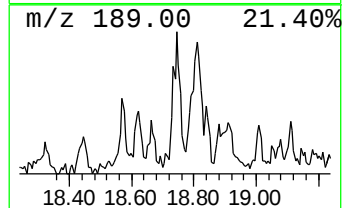
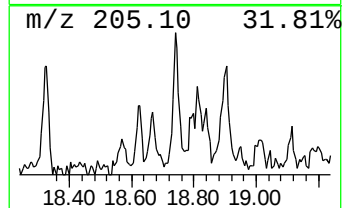
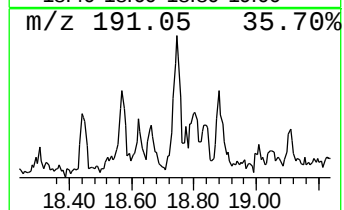
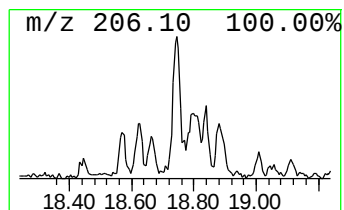
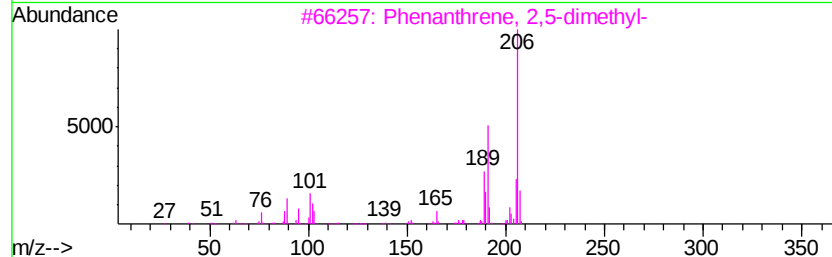
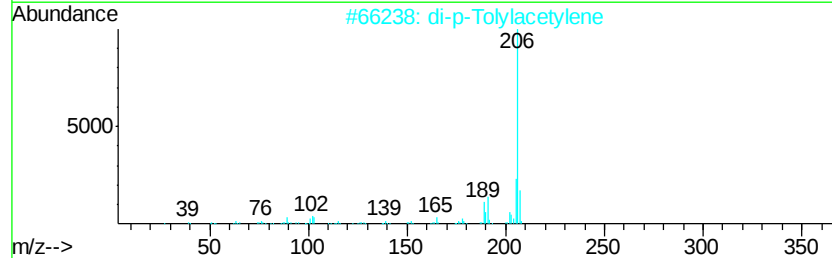
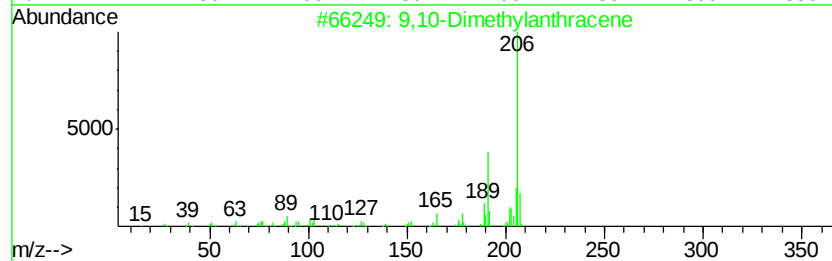
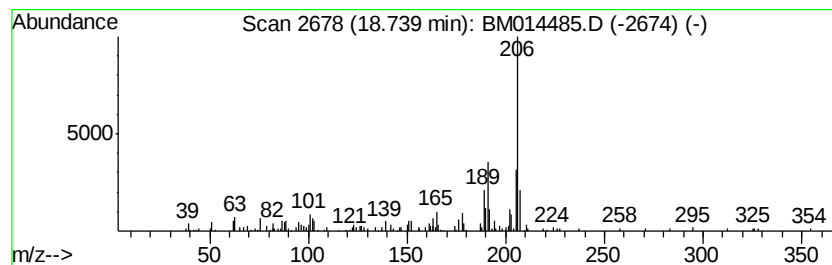
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 9,10-Dimethylantracene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.74	3.68 ng/ul	222360	Phenanthrene-d10	16.94

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	9,10-Dimethylantracene	206	C16H14	000781-43-1	91		
2	di-p-Tolylacetylene	206	C16H14	002789-88-0	91		
3	Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	91		
4	Anthracene, 1,4-dimethyl-	206	C16H14	000781-92-0	91		
5	Phenanthrene, 3,6-dimethyl-	206	C16H14	001576-67-6	91		



Data Path : Z:\HPCHEM1\BNA M\DATA\BM030518\
Data File : BM014485.D
Acq On : 05 Mar 2018 22:35
Operator : SJ/JU
Sample : J1678-09DL 5X
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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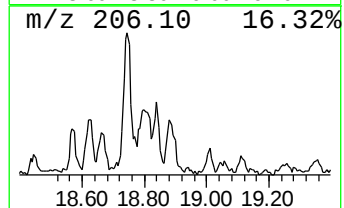
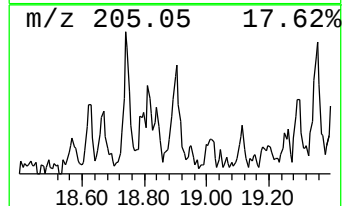
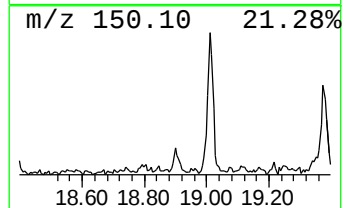
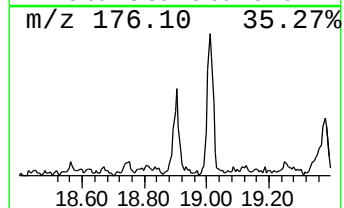
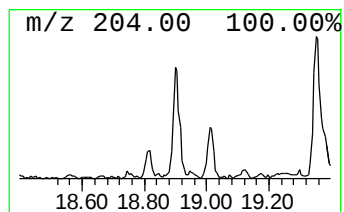
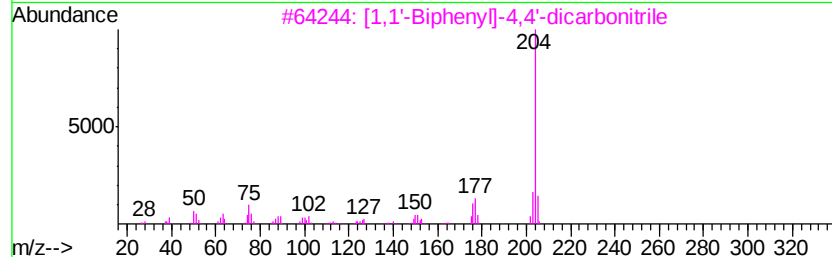
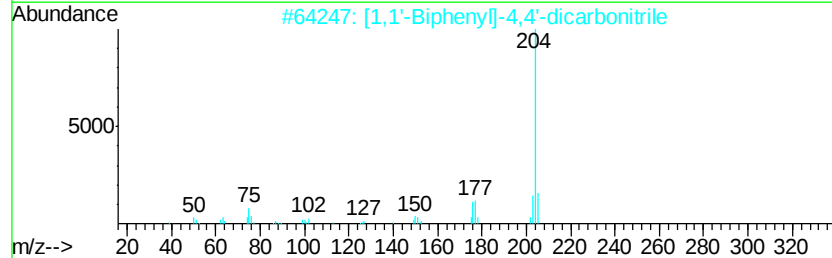
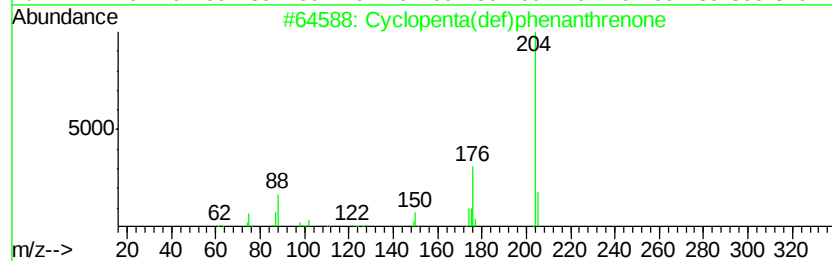
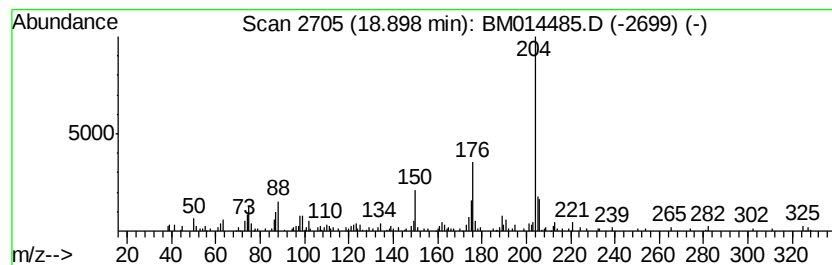
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 Cyclopenta(def)phenanthrenone Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.90	4.50 ng/ul	271941	Phenanthrene-d10	16.94

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopenta(def)phenanthrenone	204	C15H8O	005737-13-3	81
2		[1,1'-Biphenyl]-4,4'-dicarbonitrile	204	C14H8N2	001591-30-6	64
3		[1,1'-Biphenyl]-4,4'-dicarbonitrile	204	C14H8N2	001591-30-6	58
4		1,2,4,8-Tetramethylbicyclo[6.3.0...	204	C15H24	137235-51-9	55
5		Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	50



Data Path : Z:\HPCHEM1\BNA M\DATA\BM030518\
Data File : BM014485.D
Acq On : 05 Mar 2018 22:35
Operator : SJ/JU
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Misc :
ALS Vial : 18 Sample Multiplier: 1

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ClientSampleId :
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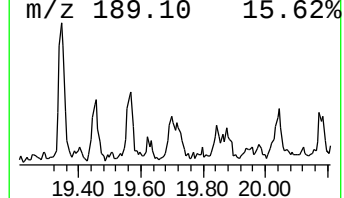
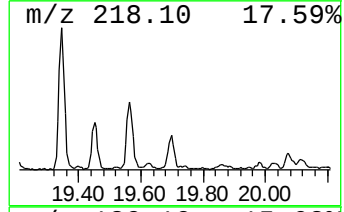
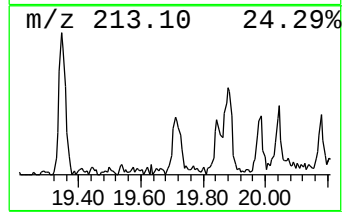
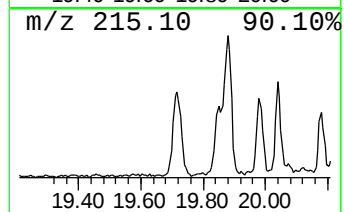
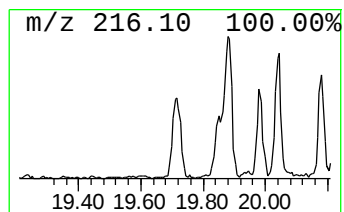
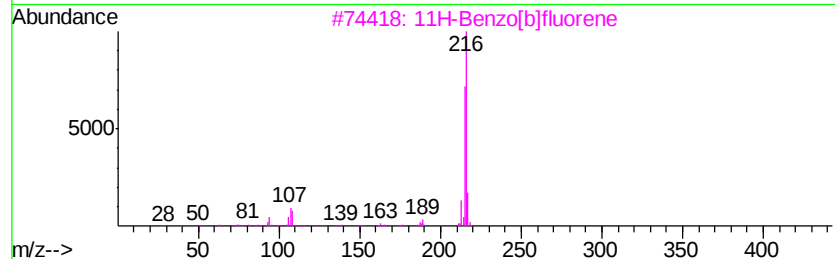
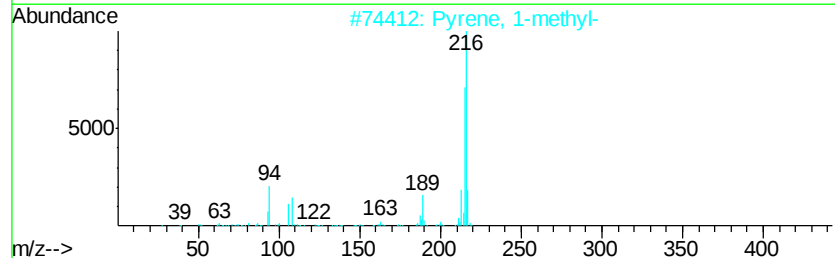
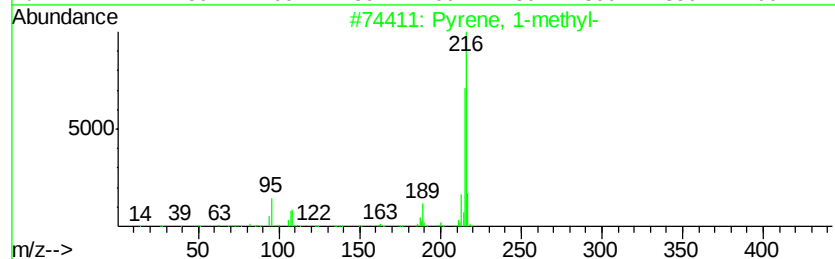
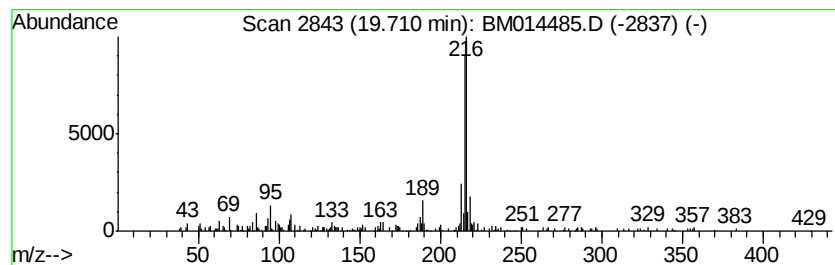
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.71	2.89 ng/ul	364692	Chrysene-d12	21.16

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pyrene, 1-methyl-	216	C17H12	002381-21-7	93
2		Pyrene, 1-methyl-	216	C17H12	002381-21-7	81
3		11H-Benzo[b]fluorene	216	C17H12	000243-17-4	78
4		Pyrene, 1-methyl-	216	C17H12	002381-21-7	74
5		11H-Benzo[b]fluorene	216	C17H12	000243-17-4	74



Data Path : Z:\HPCHEM1\BNA M\DATA\BM030518\
Data File : BM014485.D
Acq On : 05 Mar 2018 22:35
Operator : SJ/JU
Sample : J1678-09DL 5X
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
BE5T9DL

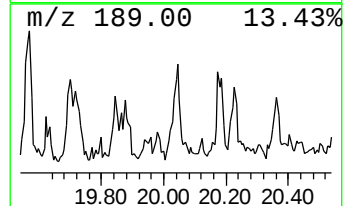
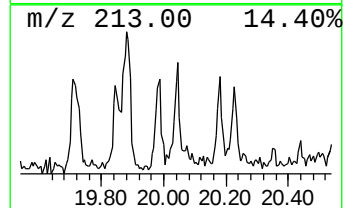
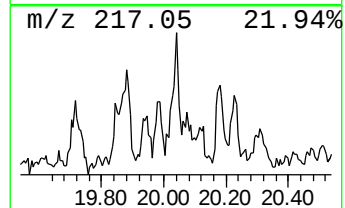
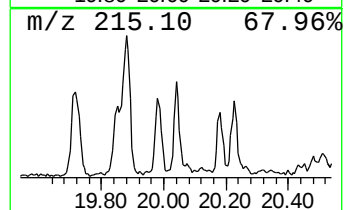
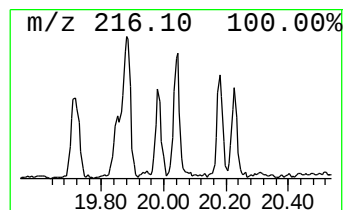
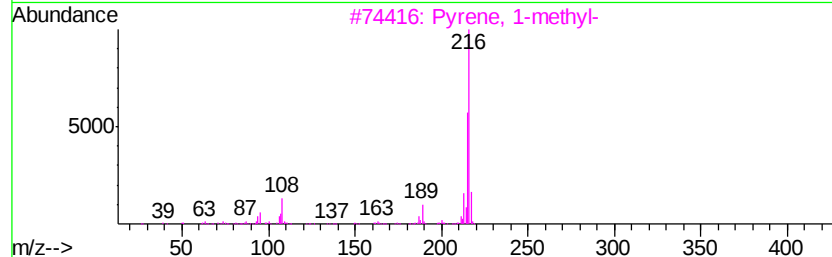
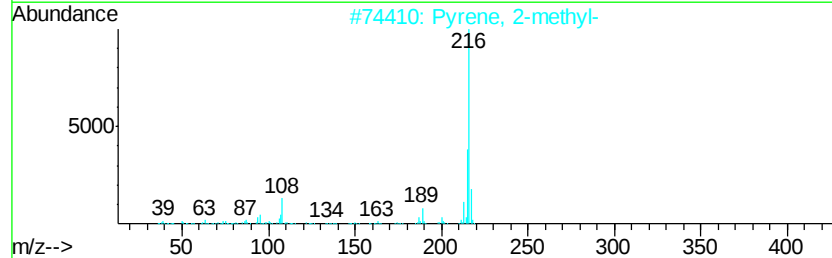
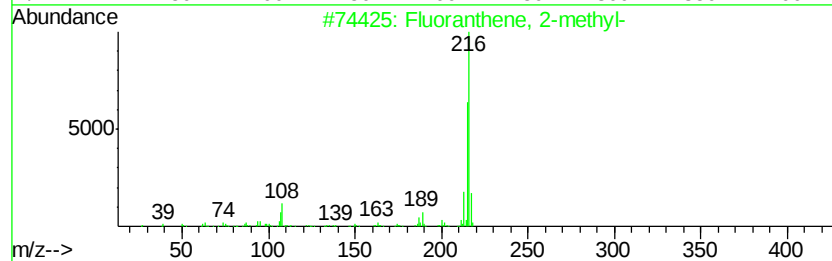
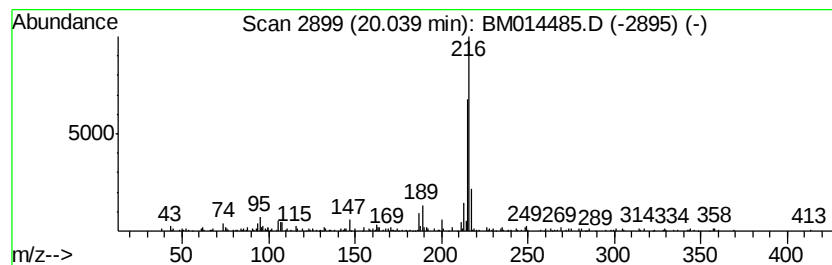
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Quant Title : SVOA CALIBRATION

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

Peak Number 13 Fluoranthene, 2-methyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.04	2.20 ng/ul	277653	Chrysene-d12	21.16

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	93
2		Pyrene, 2-methyl-	216	C17H12	003442-78-2	93
3		Pyrene, 1-methyl-	216	C17H12	002381-21-7	93
4		11H-Benzo[alfluorene	216	C17H12	000238-84-6	90
5		Pyrene, 1-methyl-	216	C17H12	002381-21-7	87



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM030518\
Data File : BM014485.D
Acq On : 05 Mar 2018 22:35
Operator : SJ/JU
Sample : J1678-09DL 5X
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
BE5T9DL

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM021418.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...	17.82	5.3	ng/ul	323016	4	16.94	1208440	20.0
Anthracene, 1-met...	17.86	6.5	ng/ul	395809	4	16.94	1208440	20.0
4H-Cyclopenta[def...	18.01	7.5	ng/ul	452229	4	16.94	1208440	20.0
Naphtho[2,3-b]nor...	18.04	3.3	ng/ul	197628	4	16.94	1208440	20.0
5,16[1',2']:8,13[...	18.32	2.8	ng/ul	168507	4	16.94	1208440	20.0
9,10-Anthracenedione	18.39	3.5	ng/ul	210260	4	16.94	1208440	20.0
Phenanthrene, 2,3...	18.57	2.6	ng/ul	156670	4	16.94	1208440	20.0
9,10-Dimethylanth...	18.74	3.7	ng/ul	222360	4	16.94	1208440	20.0
Cyclopenta(def)ph...	18.90	4.5	ng/ul	271941	4	16.94	1208440	20.0
Pyrene, 1-methyl-	19.71	2.9	ng/ul	364692	5	21.16	2520840	20.0
Fluoranthene, 2-m...	20.04	2.2	ng/ul	277653	5	21.16	2520840	20.0