

Data Path : Z:\HPCHEM1\BNA M\DATA\BM110916\
 Data File : BM007805.D
 Acq On : 09 Nov 2016 19:54
 Operator : UM/SJ
 Sample : H5554-18
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 E5N90

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\SOM02.2-EPA-BM102016.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	2.640	5	9	24	rVB4	45123	132288	2.20%	0.161%
2	6.922	728	737	752	rVB2	484313	1067035	17.73%	1.301%
3	7.081	759	764	775	rVB	395955	630486	10.48%	0.769%
4	7.281	791	798	808	rBV	516190	855930	14.22%	1.044%
5	7.739	869	876	886	rBV	746022	1260467	20.95%	1.537%
6	8.445	990	996	1010	rBV	440478	772486	12.84%	0.942%
7	8.569	1011	1017	1027	rVB	68259	121965	2.03%	0.149%
8	8.898	1066	1073	1082	rBV	486745	823388	13.68%	1.004%
9	9.616	1188	1195	1206	rBV	387316	670643	11.14%	0.818%
10	10.151	1279	1286	1304	rBV	536464	1026610	17.06%	1.252%
11	10.522	1341	1349	1356	rBV	951046	1633515	27.15%	1.992%
12	10.669	1367	1374	1390	rBV	260866	565052	9.39%	0.689%
13	13.786	1899	1904	1909	rBV	1028388	1451178	24.12%	1.770%
14	13.839	1909	1913	1921	rVB	208662	307265	5.11%	0.375%
15	14.068	1945	1952	1965	rBV	1166782	1796820	29.86%	2.192%
16	14.374	1995	2004	2011	rBV	1421357	2052929	34.12%	2.504%
17	14.439	2011	2015	2024	rVB	122852	184289	3.06%	0.225%
18	14.604	2036	2043	2053	rBV2	214509	468312	7.78%	0.571%
19	14.780	2068	2073	2079	rBV	81138	125033	2.08%	0.153%
20	15.221	2145	2148	2153	rVB	51969	65236	1.08%	0.080%
21	15.368	2167	2173	2179	rBV	1504487	2220999	36.91%	2.709%
22	15.427	2179	2183	2192	rVB	178018	259636	4.31%	0.317%
23	15.509	2192	2197	2207	rBV	292145	496185	8.25%	0.605%
24	15.621	2207	2216	2222	rVB5	24649	80072	1.33%	0.098%
25	15.727	2229	2234	2239	rBV	42162	64543	1.07%	0.079%
26	15.874	2249	2259	2266	rBV	41186	92529	1.54%	0.113%
27	16.086	2290	2295	2298	rBV	76467	117602	1.95%	0.143%
28	16.433	2344	2354	2359	rBV4	38374	93226	1.55%	0.114%
29	16.786	2409	2414	2419	rVB	87693	133886	2.22%	0.163%
30	16.874	2421	2429	2433	rBV2	85371	141577	2.35%	0.173%
31	16.945	2433	2441	2450	rVB	121207	224447	3.73%	0.274%
32	17.021	2450	2454	2461	rBV3	84798	133915	2.23%	0.163%
33	17.121	2461	2471	2475	rVV	1593429	2459068	40.87%	2.999%
34	17.168	2475	2479	2483	rVV	2491070	3495074	58.08%	4.263%

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35	17.221	2483	2488	2492	rVV2	1670019	2515111	41.80%	3.068%
36	17.256	2492	2494	2500	rVB	402946	478004	7.94%	0.583%
37	17.533	2533	2541	2551	rVB2	176821	313309	5.21%	0.382%
38	17.645	2557	2560	2566	rVB2	51037	73022	1.21%	0.089%
39	17.709	2566	2571	2575	rBV5	35269	63725	1.06%	0.078%
40	18.003	2615	2621	2625	rBV2	224688	471538	7.84%	0.575%
41	18.045	2625	2628	2633	rVB	285449	392063	6.52%	0.478%
42	18.127	2638	2642	2647	rVB	59611	86351	1.43%	0.105%
43	18.192	2647	2653	2657	rBV4	365522	606109	10.07%	0.739%
44	18.227	2657	2659	2668	rVB	130354	171802	2.86%	0.210%
45	18.309	2668	2673	2677	rBV5	51469	85033	1.41%	0.104%
46	18.503	2701	2706	2710	rBV	182890	255874	4.25%	0.312%
47	18.550	2710	2714	2719	rVV	298044	409420	6.80%	0.499%
48	18.598	2719	2722	2730	rVB5	44683	79930	1.33%	0.097%
49	18.750	2743	2748	2751	rBV2	62972	91264	1.52%	0.111%
50	18.915	2771	2776	2783	rBV4	126830	393336	6.54%	0.480%
51	19.062	2798	2801	2814	rVB2	232382	461699	7.67%	0.563%
52	19.186	2814	2822	2829	rVB	4206353	5736721	95.33%	6.997%
53	19.297	2836	2841	2845	rBV3	154356	248394	4.13%	0.303%
54	19.368	2845	2853	2860	rVV	961945	1428199	23.73%	1.742%
55	19.427	2860	2863	2872	rVB2	115545	214395	3.56%	0.261%
56	19.521	2873	2879	2881	rVV2	2663155	4030057	66.97%	4.915%
57	19.550	2881	2884	2890	rVV	3192318	4240863	70.48%	5.173%
58	19.615	2890	2895	2902	rVB2	307498	497677	8.27%	0.607%
59	19.727	2910	2914	2921	rBV	172259	236507	3.93%	0.288%
60	19.792	2921	2925	2931	rVB	59866	96011	1.60%	0.117%
61	19.868	2933	2938	2945	rVV2	160207	377580	6.27%	0.461%
62	19.933	2945	2949	2956	rVV	66176	103346	1.72%	0.126%
63	20.044	2964	2968	2973	rVB	416788	604316	10.04%	0.737%
64	20.144	2980	2985	2989	rBV	310721	442951	7.36%	0.540%
65	20.203	2989	2995	2998	rVV2	215659	409288	6.80%	0.499%
66	20.239	2999	3001	3005	rVB	89031	100048	1.66%	0.122%
67	20.344	3015	3019	3022	rBV	107578	141743	2.36%	0.173%
68	20.392	3023	3027	3030	rVB2	121324	170075	2.83%	0.207%
69	20.797	3092	3096	3101	rBV	264713	352184	5.85%	0.430%
70	20.956	3118	3123	3126	rBV2	374817	528830	8.79%	0.645%
71	20.997	3126	3130	3131	rVV	288112	360786	6.00%	0.440%

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72	21.027	3131	3135	3141	rVV3	366416	926150	15.39%	1.130%
73	21.091	3142	3146	3151	rVB2	323788	454521	7.55%	0.554%
74	21.309	3176	3183	3187	rBV2	3031967	6017528	100.00%	7.339%
75	21.344	3187	3189	3199	rVB	1873653	2278247	37.86%	2.779%
76	21.462	3206	3209	3216	rVB	319403	404853	6.73%	0.494%
77	21.627	3233	3237	3242	rVB2	227318	333597	5.54%	0.407%
78	22.850	3441	3445	3447	rBV	698670	951054	15.80%	1.160%
79	22.891	3447	3452	3465	rVB	1596684	4965925	82.52%	6.057%
80	23.368	3528	3533	3537	rBV	817340	1448583	24.07%	1.767%
81	23.421	3537	3542	3546	rVV2	1659603	3017116	50.14%	3.680%
82	23.462	3546	3549	3557	rVB	1201212	2122809	35.28%	2.589%
83	23.562	3560	3566	3571	rBV	1590399	2903283	48.25%	3.541%
84	24.062	3646	3651	3656	rVB2	390254	710432	11.81%	0.867%
85	25.826	3945	3951	3963	rVB2	755382	2163299	35.95%	2.639%

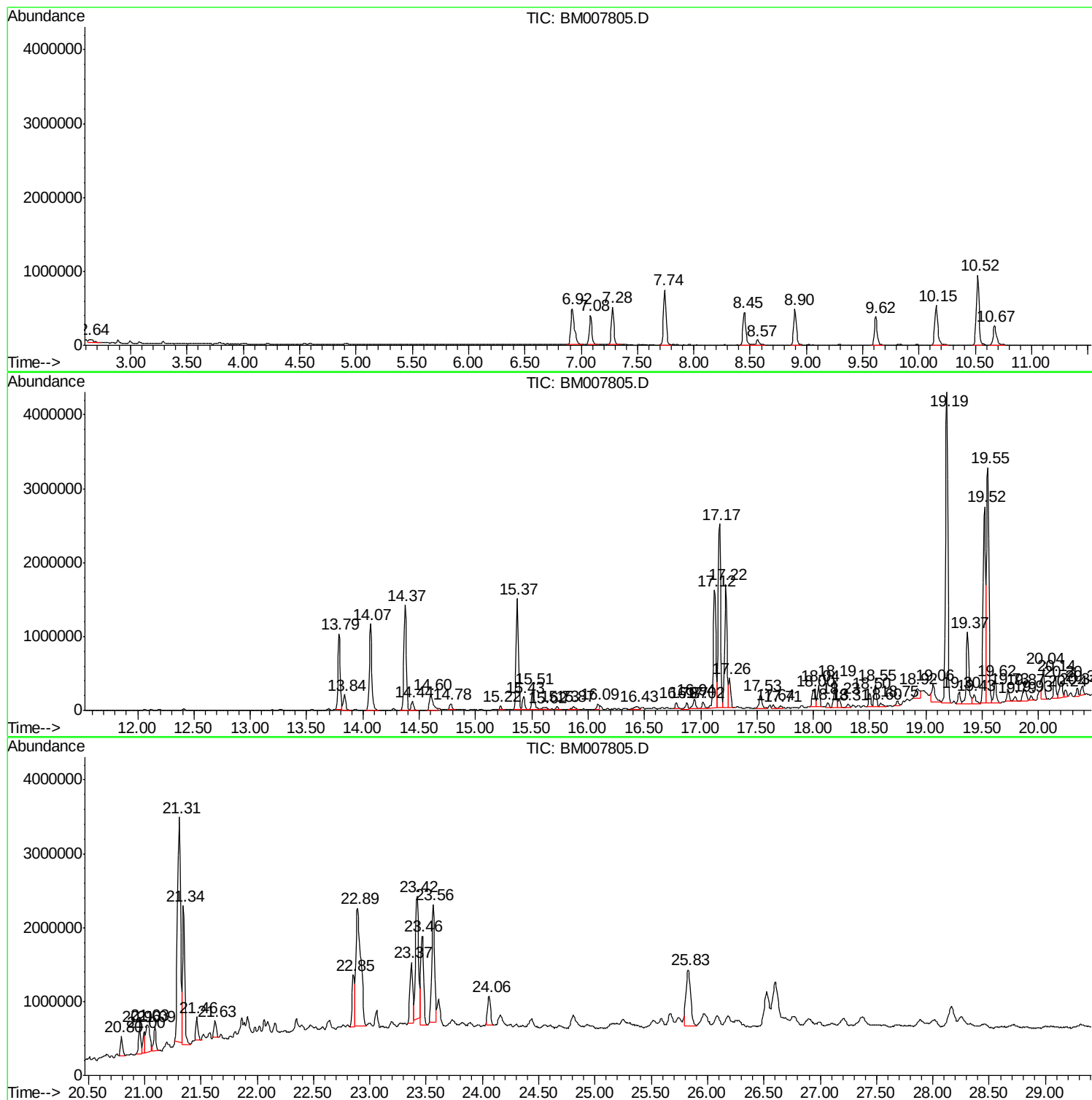
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ClientSampled :
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Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM102016.M
Quant Title : SVOA CALIBRATION

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



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Instrument :
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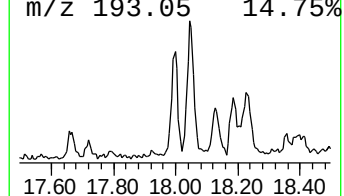
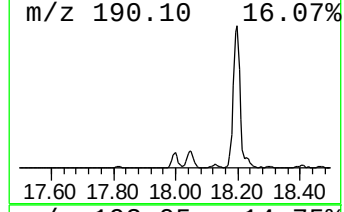
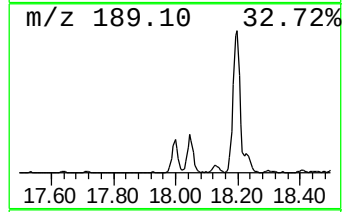
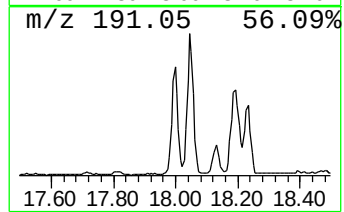
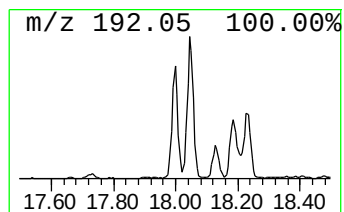
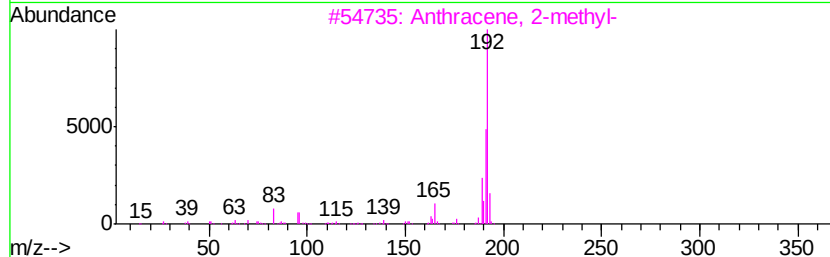
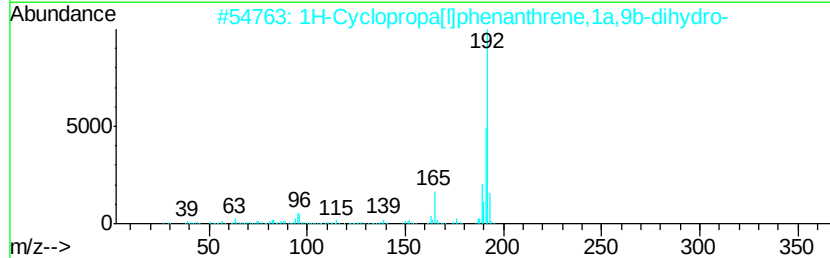
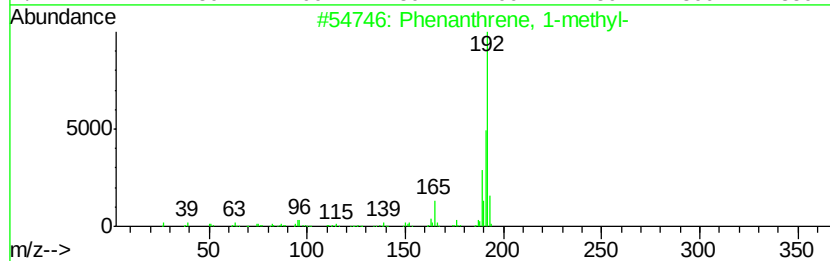
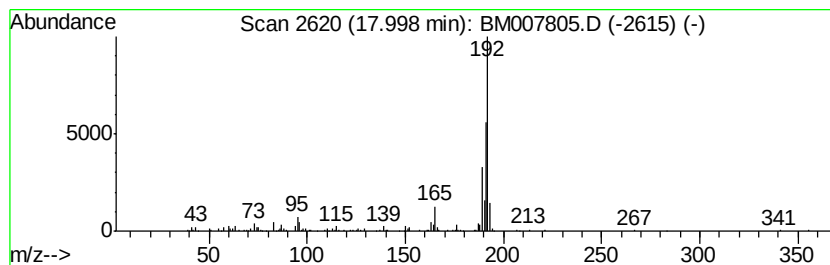
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM102016.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Phenanthrene, 1-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.00	3.84 ng/ul	471538	Phenanthrene-d10	17.12

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	96
2		1H-Cyclopropa[1]phenanthrene,1a,...	192	C15H12	000949-41-7	95
3		Anthracene, 2-methyl-	192	C15H12	000613-12-7	89
4		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	89
5		Anthracene, 9-methyl-	192	C15H12	000779-02-2	89



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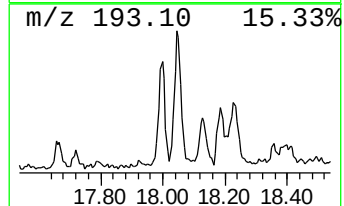
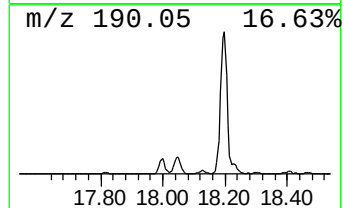
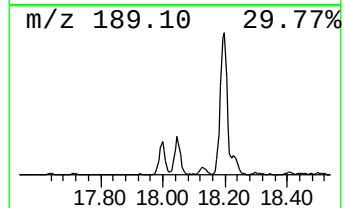
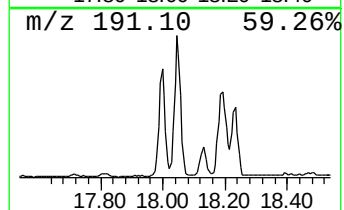
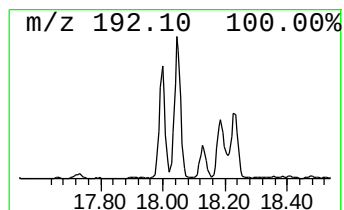
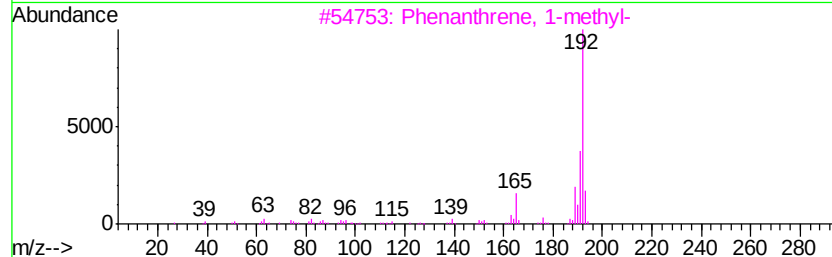
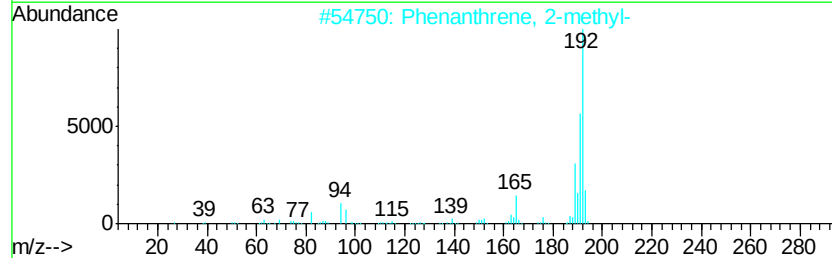
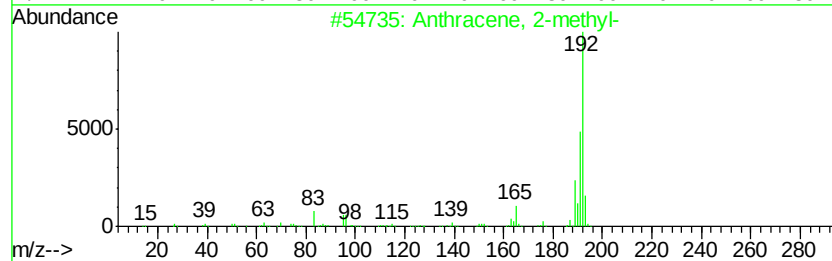
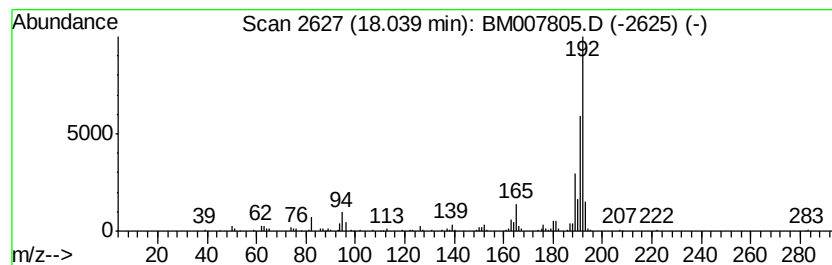
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM102016.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 Anthracene, 2-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.04	3.19 ng/ul	392063	Phenanthrene-d10	17.12

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Anthracene, 2-methyl-	192	C15H12	000613-12-7	95
2		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	95
3		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	95
4		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	94
5		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	94



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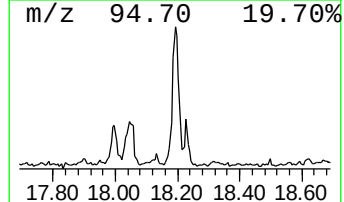
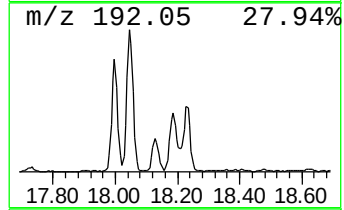
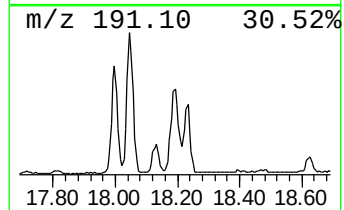
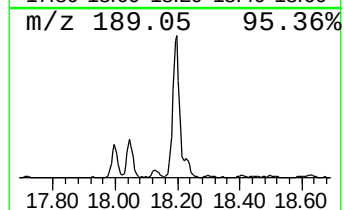
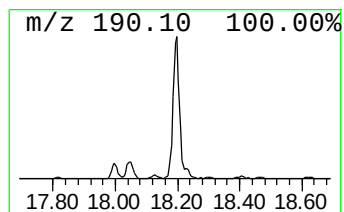
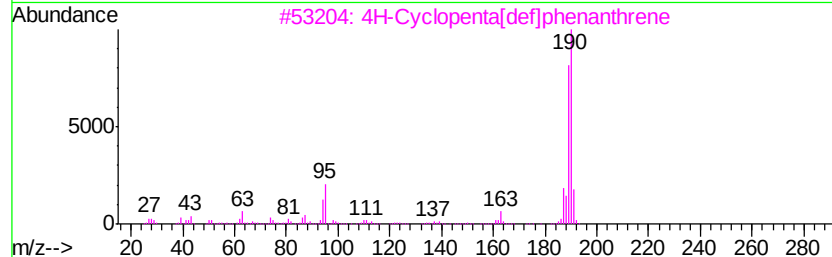
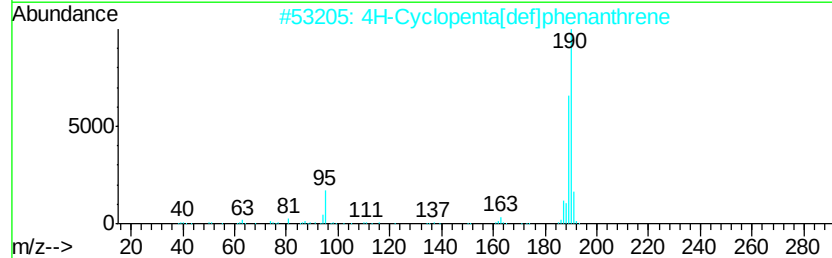
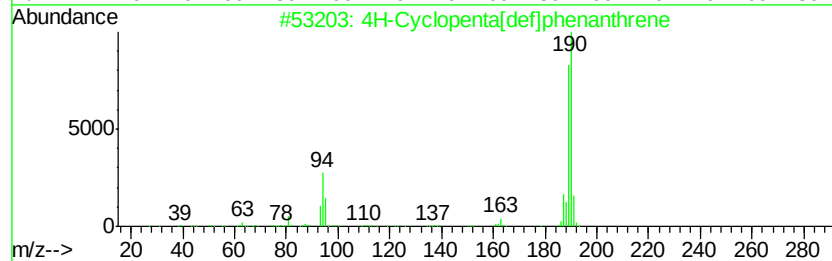
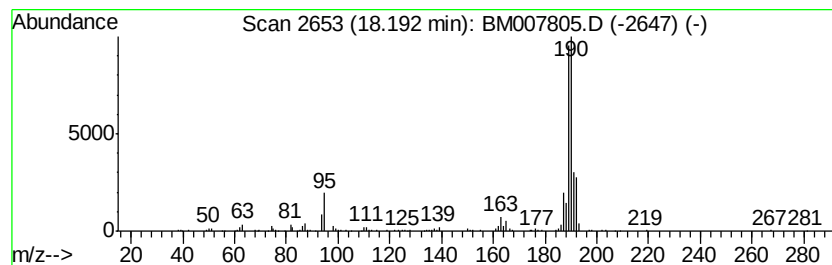
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Quant Title : SVOA CALIBRATION

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

R.T.	EstConc	Area	Relative to ISTD	R.T.		
18.19	4.93 ng/ul	606109	Phenanthrene-d10	17.12		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	87
2		4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	81
3		4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	60
4		Pyrazole-4-carboxaldehyde, 3-(4-...	190	C10H7FN2O	306936-57-2	53
5		2,3,5,6-Tetramethylterephthalald...	190	C12H14O2	007072-01-7	50



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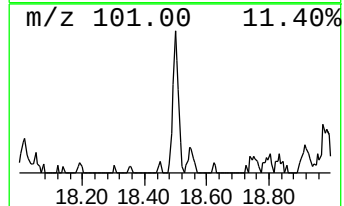
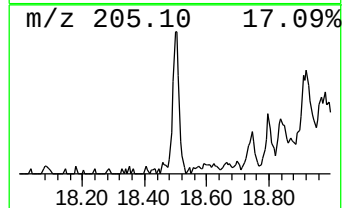
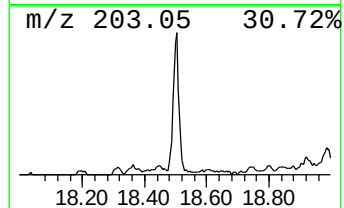
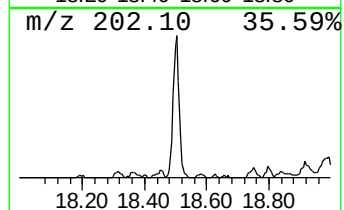
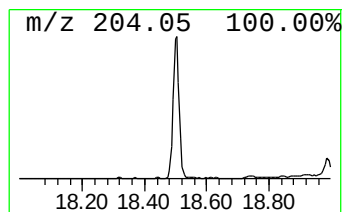
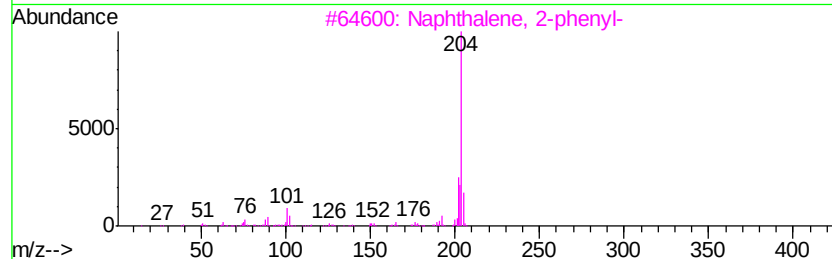
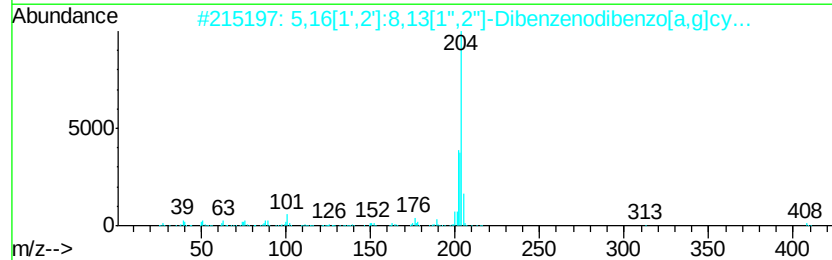
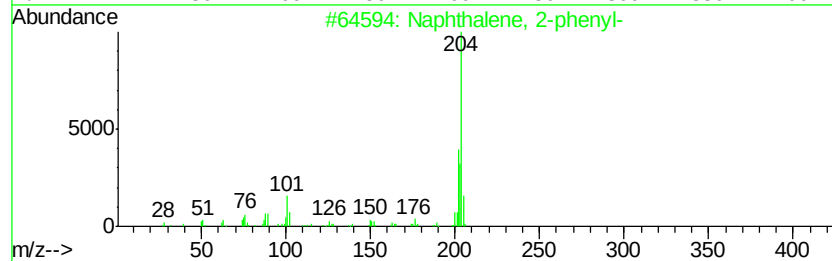
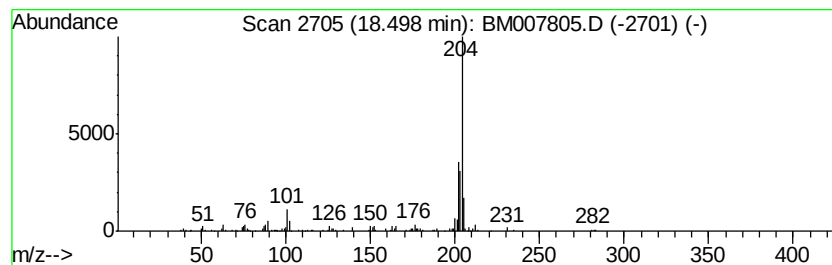
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Naphthalene, 2-phenyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.50	2.08 ng/ul	255874	Phenanthrene-d10	17.12

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	93
2		5,16[1',2']:8,13[1'',2'']-Dibenz...	408	C32H24	005672-97-9	90
3		Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	89
4		Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	81
5		Tricyclo[8.2.2.2(4,7)]hexadeca-2...	204	C16H12	006572-60-7	74



Data Path : Z:\HPCHEM1\BNA M\DATA\BM110916\
Data File : BM007805.D
Acq On : 09 Nov 2016 19:54
Operator : UM/SJ
Sample : H5554-18
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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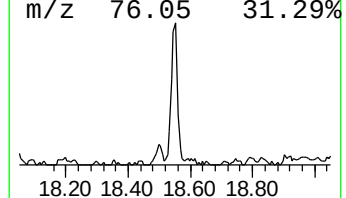
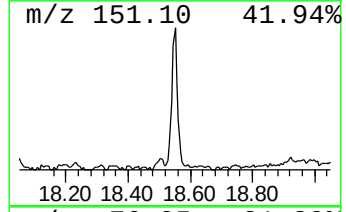
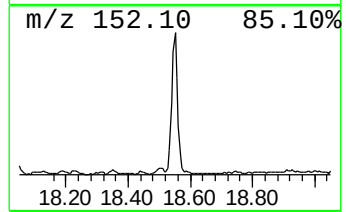
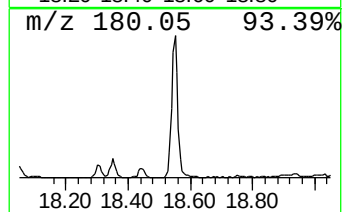
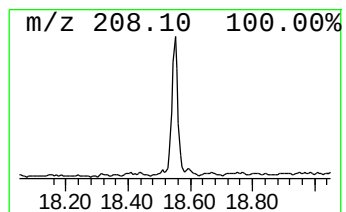
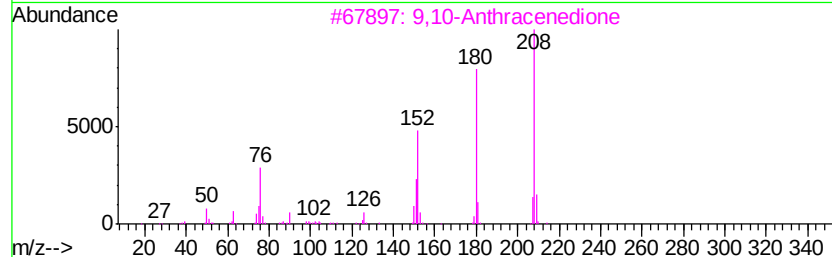
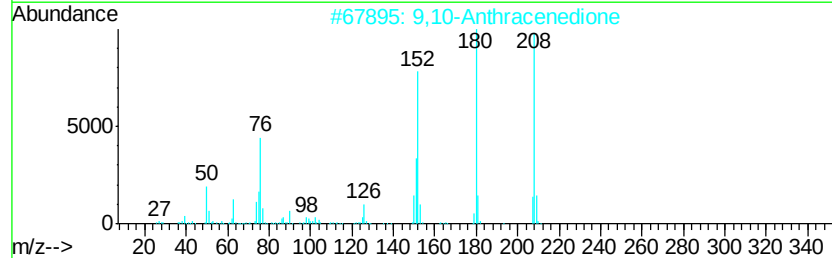
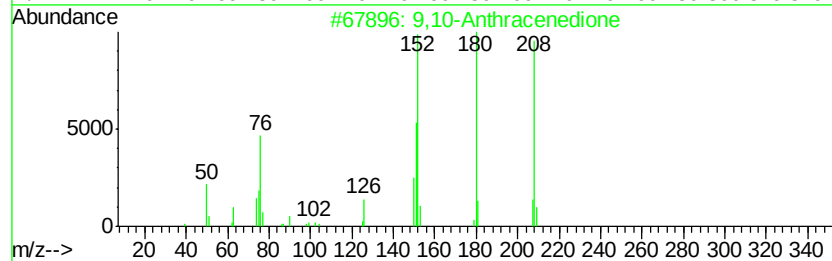
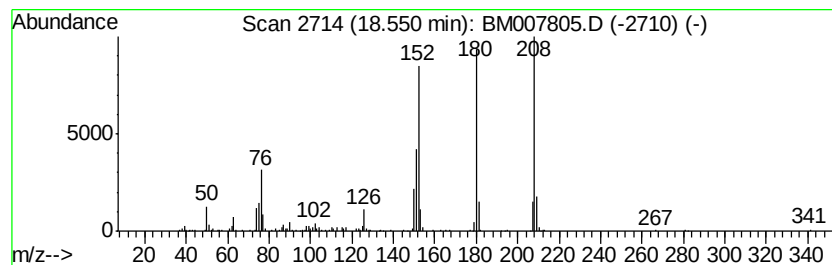
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 9,10-Anthracenedione Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.55	3.33 ng/ul	409420	Phenanthrene-d10	17.12

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	9,10-Anthracenedione	208	C14H8O2	000084-65-1	98
2		9,10-Anthracenedione	208	C14H8O2	000084-65-1	95
3		9,10-Anthracenedione	208	C14H8O2	000084-65-1	94
4		Benzo[c]cinnoline	180	C12H8N2	000230-17-1	86
5		Benzo[c]cinnoline	180	C12H8N2	000230-17-1	70



Data Path : Z:\HPCHEM1\BNA M\DATA\BM110916\
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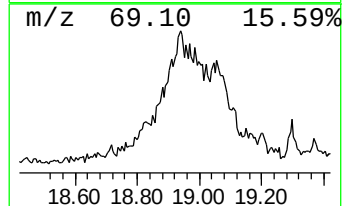
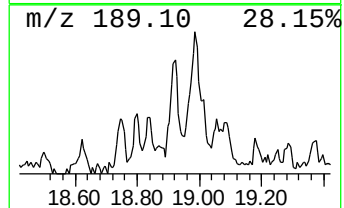
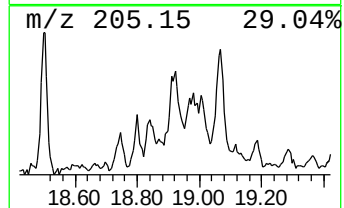
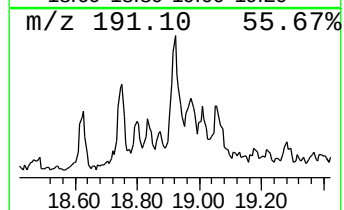
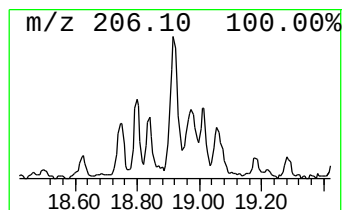
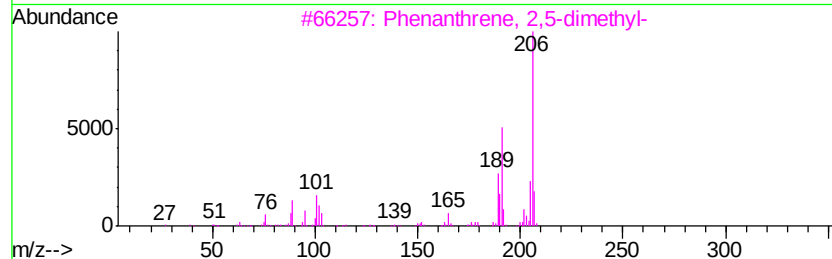
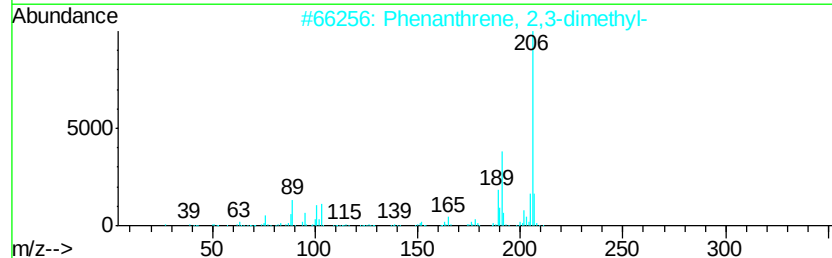
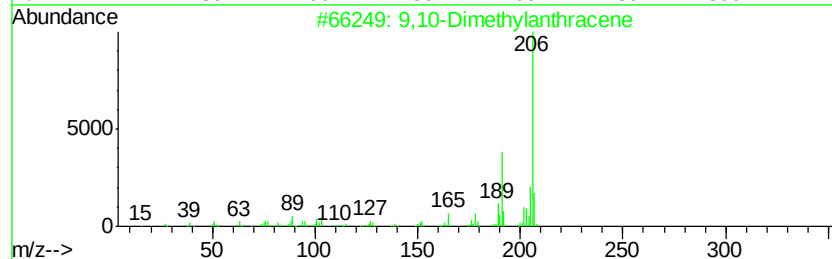
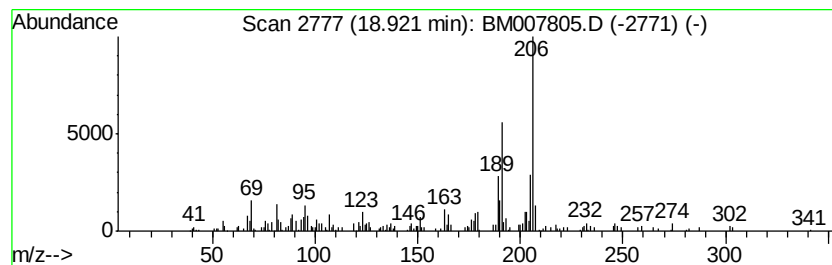
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 9,10-Dimethylantracene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.92	3.20 ng/ul	393336	Phenanthrene-d10	17.12

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			9,10-Dimethylantracene	206	C16H14	000781-43-1	95
2			Phenanthrene, 2,3-dimethyl-	206	C16H14	003674-65-5	93
3			Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	90
4			Anthracene, 1,4-dimethyl-	206	C16H14	000781-92-0	86
5			Phenanthrene, 3,6-dimethyl-	206	C16H14	001576-67-6	81



Data Path : Z:\HPCHEM1\BNA M\DATA\BM110916\
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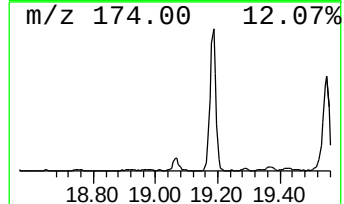
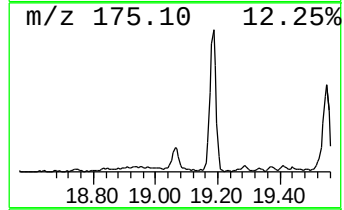
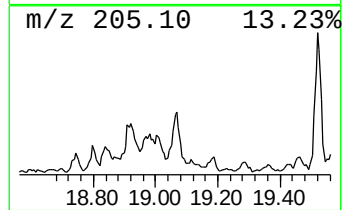
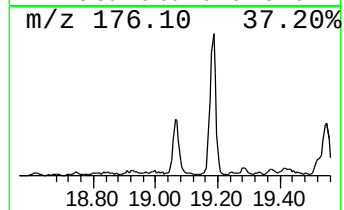
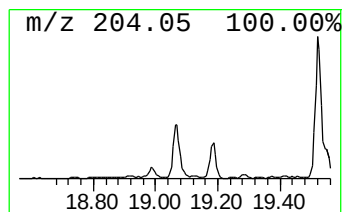
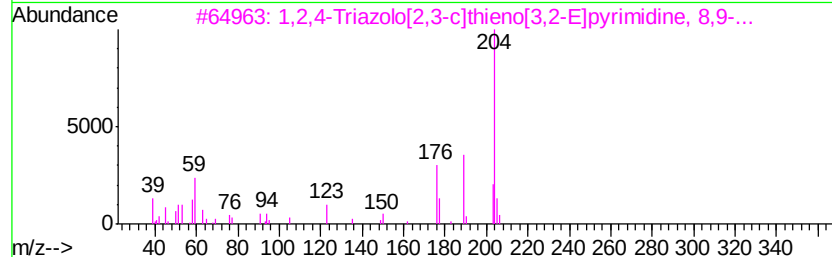
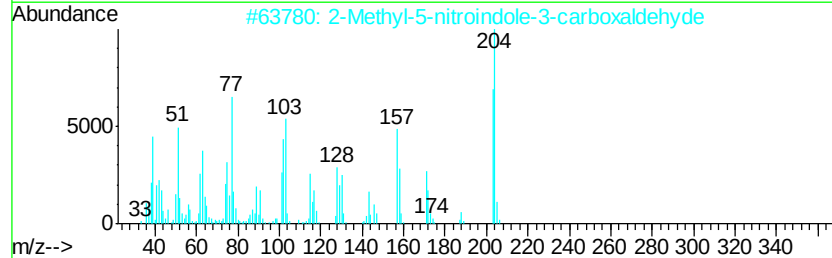
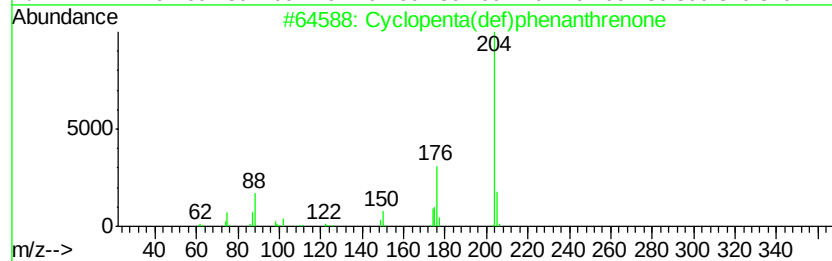
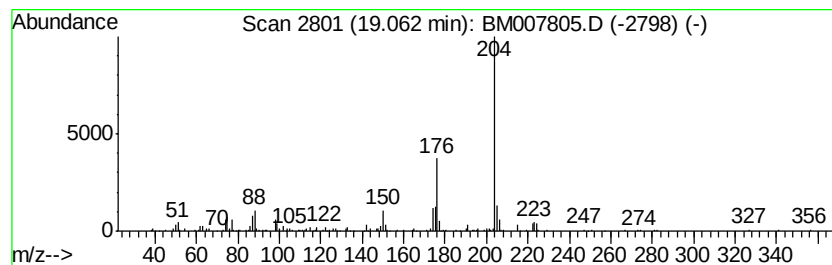
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Cyclopenta(def)phenanthrenone Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.06	3.76 ng/ul	461699	Phenanthrene-d10	17.12

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopenta(def)phenanthrenone	204	C15H8O	005737-13-3	72
2		2-Methyl-5-nitroindole-3-carboxal...	204	C10H8N2O3	003558-17-6	46
3		1,2,4-Triazolo[2,3-c]thieno[3,2-...	204	C9H8N4S	113246-88-1	45
4		6-Phenyl-2-thiouracil	204	C10H8N2OS	005590-94-3	45
5		[1,1'-Biphenyl]-4,4'-dicarbonitrile	204	C14H8N2	001591-30-6	45



Data Path : Z:\HPCHEM1\BNA M\DATA\BM110916\
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Sample : H5554-18
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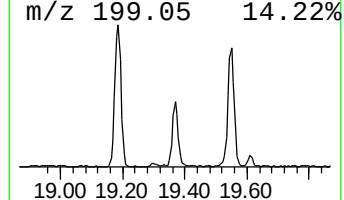
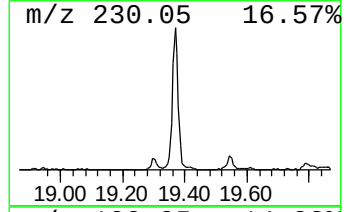
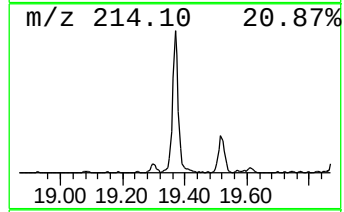
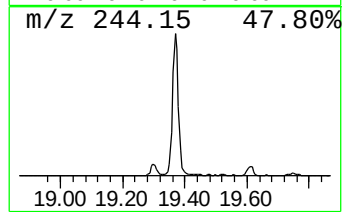
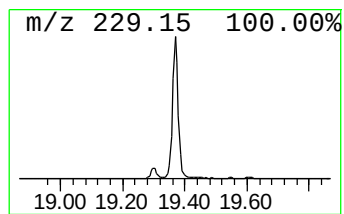
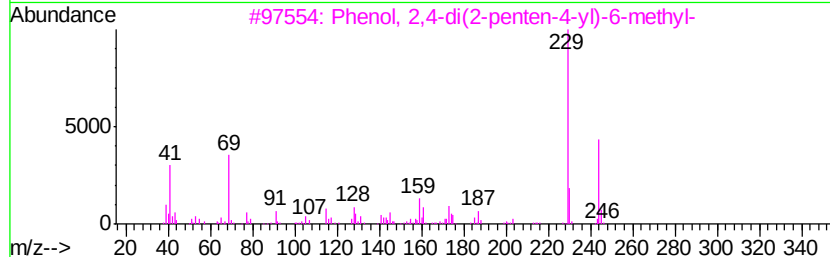
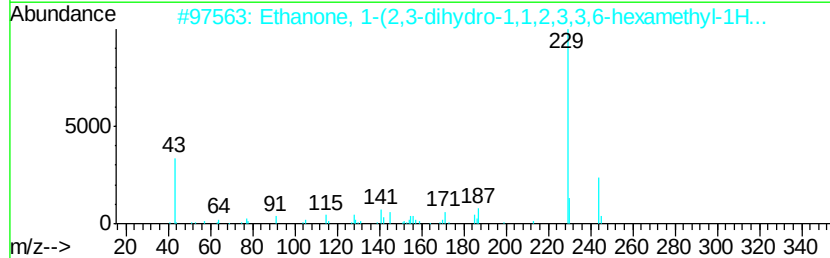
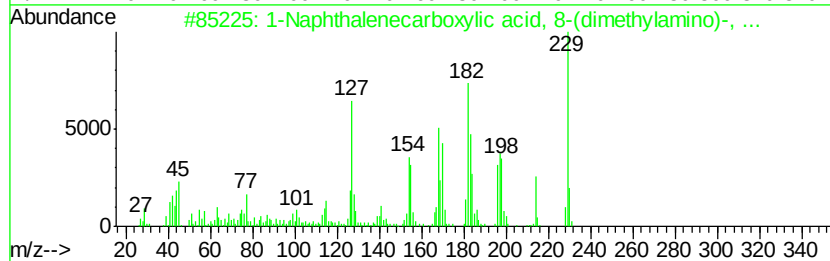
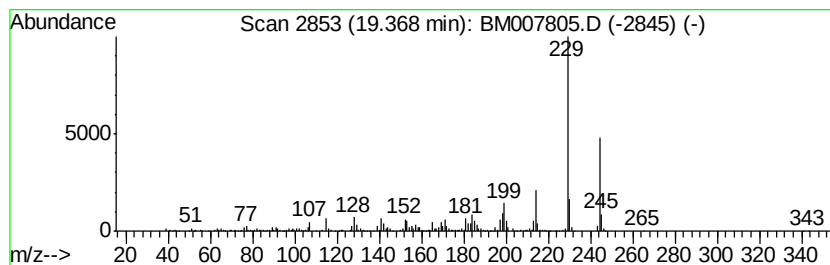
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 1-Naphthalenecarboxylic aci... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.37	4.75 ng/ul	1428200	Chrysene-d12	21.31

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Naphthalenecarboxylic acid, 8-...	229	C14H15N02	069674-55-1	86
2		Ethanone, 1-(2,3-dihydro-1,1,2,3...	244	C17H24O	015323-35-0	70
3		Phenol, 2,4-di(2-penten-4-yl)-6-...	244	C17H24O	017258-43-4	70
4		6-Acetyl-5-hydroxy-2,7-dimethyl-...	244	C14H12O4	001086-07-3	68
5		2-Propenoic acid, 2-cyano-3-[4-(...	244	C14H16N2O2	1000115-72-8	68



Data Path : Z:\HPCHEM1\BNA M\DATA\BM110916\
Data File : BM007805.D
Acq On : 09 Nov 2016 19:54
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Misc :
ALS Vial : 15 Sample Multiplier: 1

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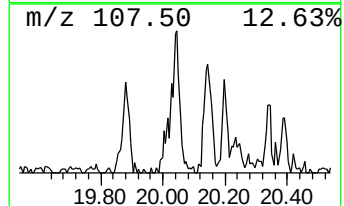
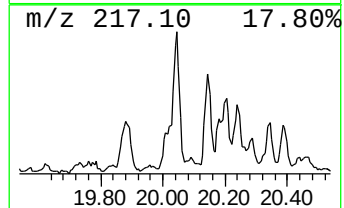
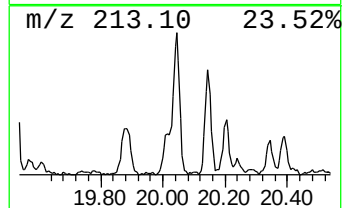
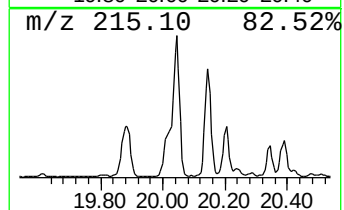
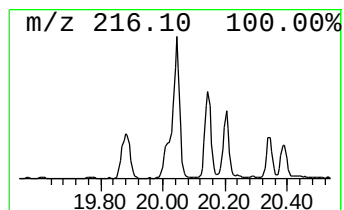
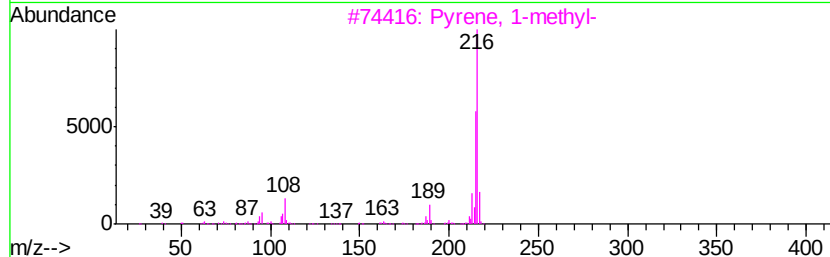
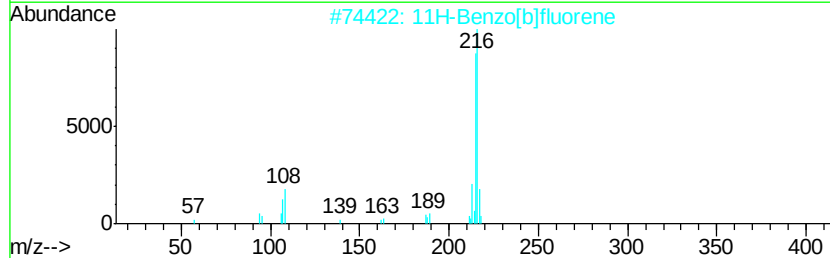
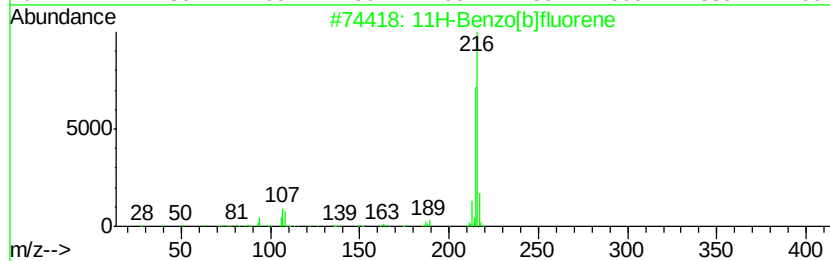
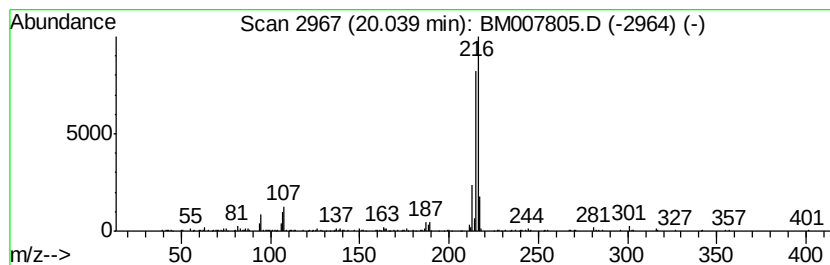
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 11H-Benzo[b]fluorene Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.04	2.01 ng/ul	604316	Chrysene-d12	21.31

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	93
2		11H-Benzo[b]fluorene	216	C17H12	000243-17-4	91
3		Pvrene, 1-methyl-	216	C17H12	002381-21-7	86
4		Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	81
5		11H-Benzo[a]fluorene	216	C17H12	000238-84-6	81



Data Path : Z:\HPCHEM1\BNA M\DATA\BM110916\
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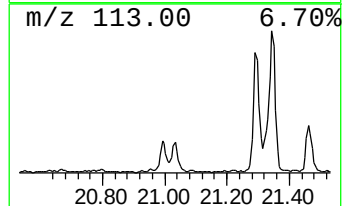
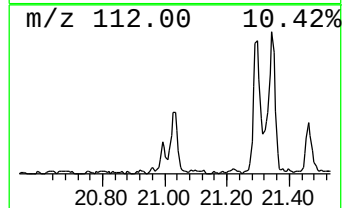
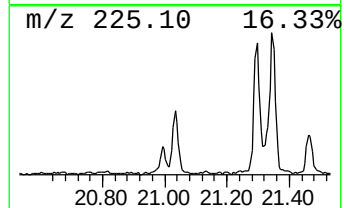
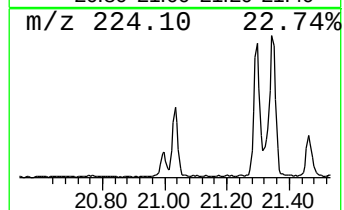
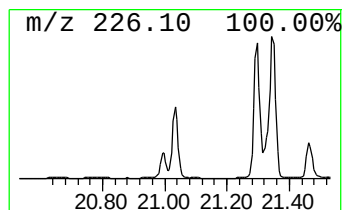
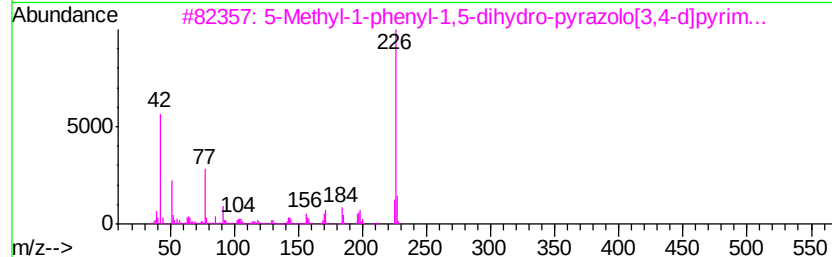
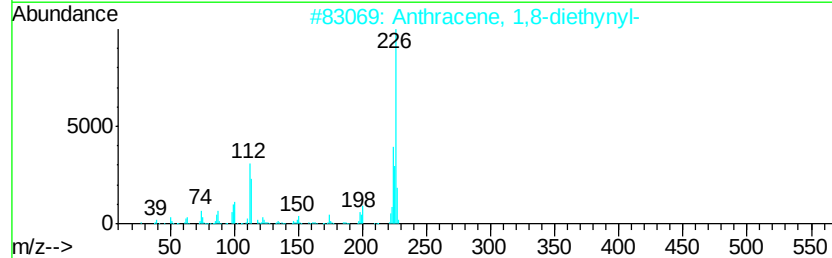
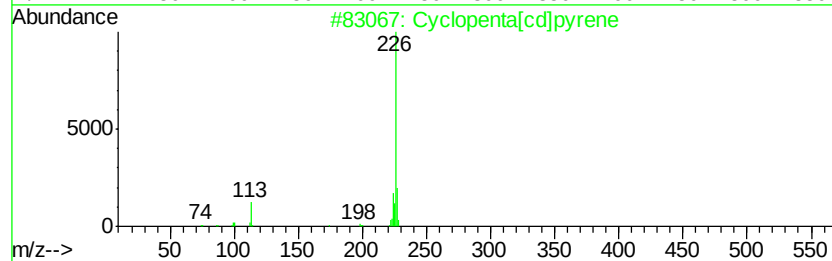
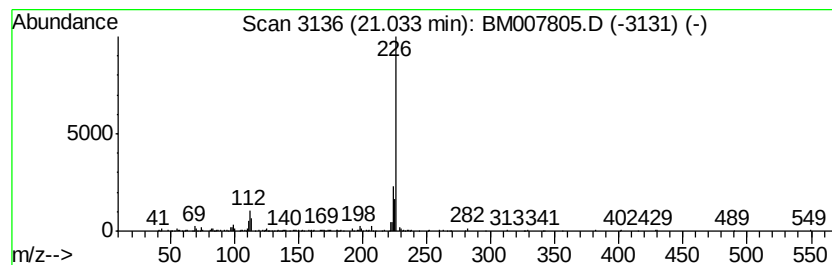
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 unknown-01 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.03	3.08 ng/ul	926150	Chrysene-d12	21.31

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopenta[cd]pyrene	226	C18H10	027208-37-3	47
2		Anthracene, 1,8-diethynyl-	226	C18H10	078053-58-4	38
3		5-Methyl-1-phenyl-1,5-dihydro-py...	226	C12H10N4O	1000300-02-1	28
4		Benzo[ghi]fluoranthene	226	C18H10	000203-12-3	25
5		N-2-Furoylglycine, tert-butyldim...	283	C13H21N04Si	1000332-85-8	23



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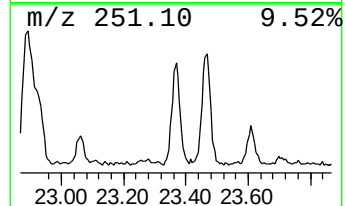
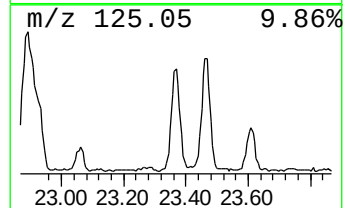
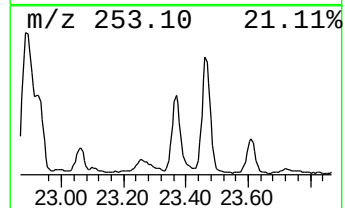
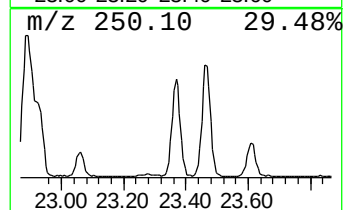
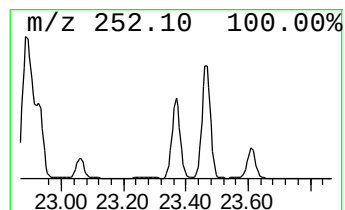
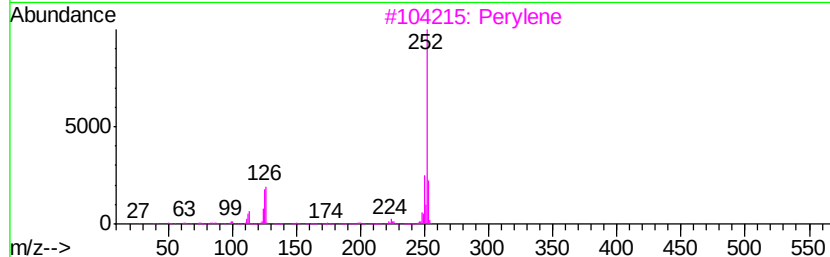
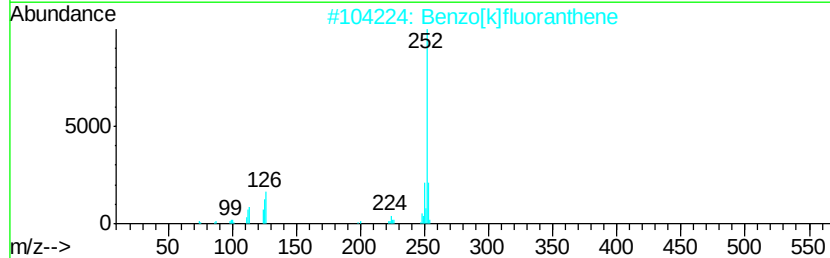
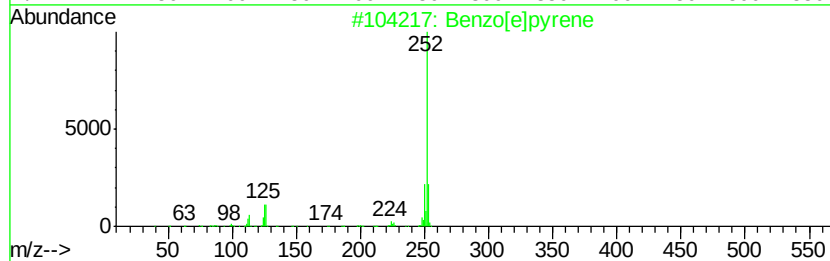
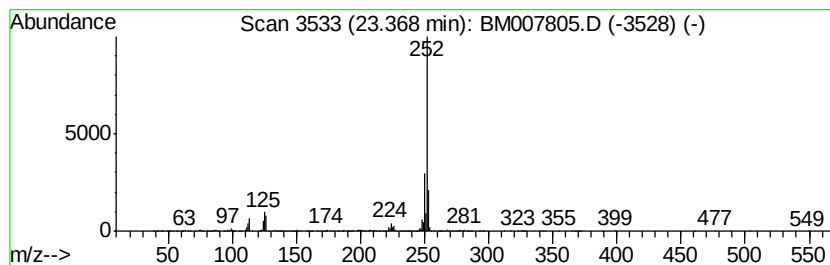
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 Benzo[e]pyrene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.37	9.98 ng/ul	1448580	Perylene-d12	23.56

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzo[e]pyrene	252	C20H12	000192-97-2	97
2		Benzo[k]fluoranthene	252	C20H12	000207-08-9	97
3		Perylene	252	C20H12	000198-55-0	94
4		Perylene	252	C20H12	000198-55-0	93
5		Benz[e]acephenanthrylene	252	C20H12	000205-99-2	91



Data Path : Z:\HPCHEM1\BNA M\DATA\BM110916\
Data File : BM007805.D
Acq On : 09 Nov 2016 19:54
Operator : UM/SJ
Sample : H5554-18
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
E5N90

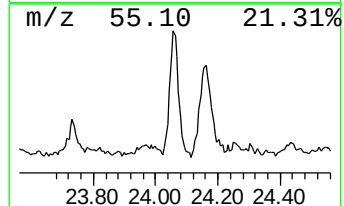
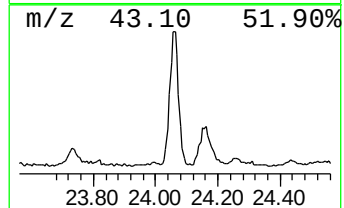
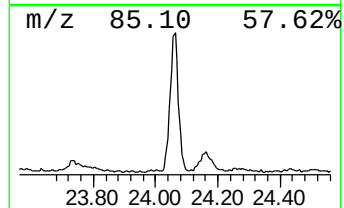
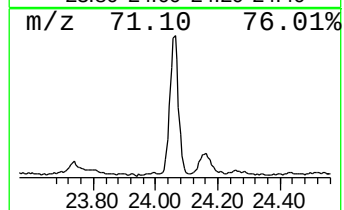
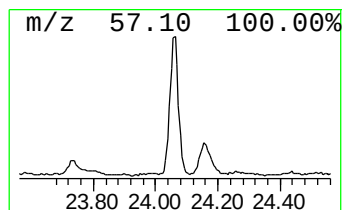
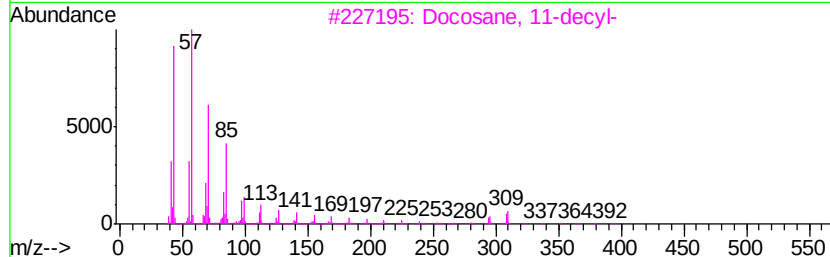
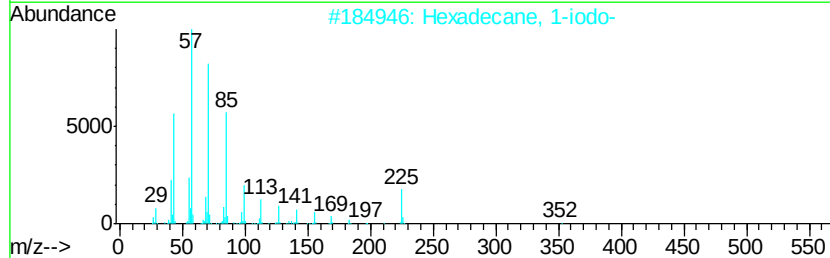
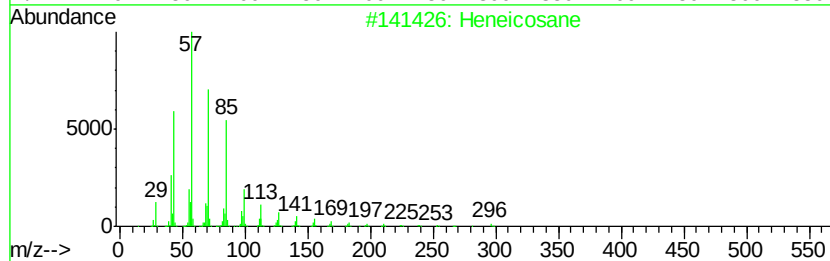
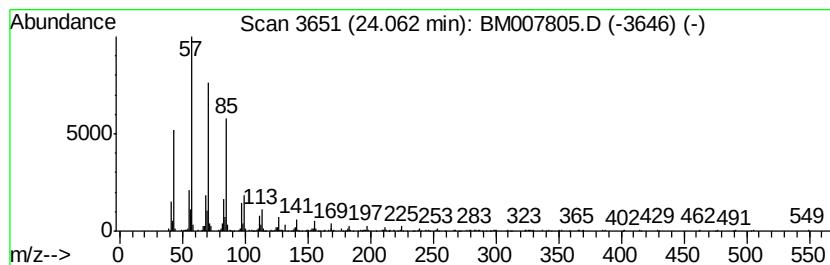
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Quant Title : SVOA CALIBRATION

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

Peak Number 13 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.06	4.89 ng/ul	710432	Perylene-d12	23.56

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	94
2		Hexadecane, 1-iodo-	352	C16H33I	000544-77-4	93
3		Docosane, 11-decyl-	451	C32H66	055401-55-3	91
4		2-methyloctacosane	408	C29H60	1000376-72-8	91
5		Triacontane	422	C30H62	000638-68-6	91



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM110916\
Data File : BM007805.D
Acq On : 09 Nov 2016 19:54
Operator : UM/SJ
Sample : H5554-18
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
E5N90

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM102016.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...	18.00	3.8	ng/ul	471538	4	17.12	2459070	20.0
Anthracene, 2-met...	18.04	3.2	ng/ul	392063	4	17.12	2459070	20.0
4H-Cyclopenta[def...	18.19	4.9	ng/ul	606109	4	17.12	2459070	20.0
Naphthalene, 2-ph...	18.50	2.1	ng/ul	255874	4	17.12	2459070	20.0
9,10-Anthracenedione	18.55	3.3	ng/ul	409420	4	17.12	2459070	20.0
9,10-Dimethylanthe...	18.92	3.2	ng/ul	393336	4	17.12	2459070	20.0
Cyclopenta(def)ph...	19.06	3.8	ng/ul	461699	4	17.12	2459070	20.0
1-Naphthalenecarb...	19.37	4.8	ng/ul	1428200	5	21.31	6017530	20.0
11H-Benzo[b]fluorene	20.04	2.0	ng/ul	604316	5	21.31	6017530	20.0
unknown-01	21.03	3.1	ng/ul	926150	5	21.31	6017530	20.0
Benzo[e]pyrene	23.37	10.0	ng/ul	1448580	6	23.56	2903280	20.0
(DEL) Alkane: Str...	24.06	4.9	ng/ul	710432	6	23.56	2903280	20.0