

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM032322\
 Data File : BM034354.D
 Acq On : 24 Mar 2022 00:21
 Operator : CG/JU
 Sample : N1958-08 5X
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C-8-7-8

Integration Parameters: rteint.p

Integrator: RTE

Smoothing : ON

Filtering: 5

Sampling : 1

Min Area: 3 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM031522.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BM034354.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.496	378	384	392	rBV	521704	784157	13.39%	1.599%
2	7.096	648	656	674	rBV	487086	834935	14.26%	1.703%
3	7.937	792	799	809	rBV	683676	1091667	18.65%	2.226%
4	9.095	990	996	1008	rBV	290179	519525	8.87%	1.059%
5	10.748	1269	1277	1283	rBV	888091	1496100	25.56%	3.051%
6	10.801	1283	1286	1292	rVB	98811	153160	2.62%	0.312%
7	12.413	1547	1560	1565	rBV2	37644	73345	1.25%	0.150%
8	12.631	1590	1597	1604	rVB	35116	58645	1.00%	0.120%
9	13.207	1688	1695	1707	rBV	770011	1153240	19.70%	2.352%
10	13.901	1807	1813	1819	rBV	33162	61205	1.05%	0.125%
11	14.307	1874	1882	1889	rBV	41102	75129	1.28%	0.153%
12	14.407	1893	1899	1908	rBV2	36059	82533	1.41%	0.168%
13	14.577	1920	1928	1934	rBV	1251767	1851070	31.62%	3.775%
14	14.642	1934	1939	1946	rVB	155845	231849	3.96%	0.473%
15	14.977	1990	1996	2003	rVB2	89923	142596	2.44%	0.291%
16	15.619	2100	2105	2111	rBV2	111590	176532	3.02%	0.360%
17	15.919	2151	2156	2164	rVB	66121	115944	1.98%	0.236%
18	16.060	2173	2180	2189	rBV	632703	965171	16.49%	1.968%
19	16.295	2215	2220	2230	rVB3	52942	114179	1.95%	0.233%
20	16.624	2267	2276	2281	rBV4	49706	118856	2.03%	0.242%
21	16.854	2309	2315	2319	rBV3	47166	86778	1.48%	0.177%
22	16.971	2331	2335	2343	rVB	85438	135227	2.31%	0.276%
23	17.054	2344	2349	2351	rVV2	49449	84208	1.44%	0.172%
24	17.083	2351	2354	2359	rVV2	61593	102748	1.76%	0.210%
25	17.136	2359	2363	2372	rVB	131836	216741	3.70%	0.442%
26	17.318	2388	2394	2398	rBV	1413012	2066004	35.29%	4.213%
27	17.360	2398	2401	2408	rVV	2029103	2791453	47.68%	5.693%
28	17.454	2412	2417	2422	rVV	421889	616403	10.53%	1.257%
29	17.513	2422	2427	2432	rVV4	46911	81362	1.39%	0.166%
30	17.718	2459	2462	2466	rBV	54591	76152	1.30%	0.155%
31	17.818	2476	2479	2486	rBV2	72345	145550	2.49%	0.297%
32	18.201	2539	2544	2548	rBV	238360	446232	7.62%	0.910%
33	18.242	2548	2551	2560	rVB	335139	500815	8.55%	1.021%
34	18.324	2561	2565	2570	rVV	134644	188166	3.21%	0.384%
35	18.389	2571	2576	2580	rVV2	345658	651321	11.13%	1.328%
36	18.424	2580	2582	2588	rVB	165437	220735	3.77%	0.450%

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Peak Location: TOP

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37	18.512	2594	2597	2601	rBV	101449	110730	1.89%	0.226%
38	18.695	2624	2628	2631	rBV	190836	249597	4.26%	0.509%
39	18.730	2631	2634	2643	rVB	180873	257909	4.41%	0.526%
40	19.112	2695	2699	2702	rBV	145867	213330	3.64%	0.435%
41	19.142	2702	2704	2708	rVV	101734	116127	1.98%	0.237%
42	19.248	2719	2722	2729	rVB	166967	253766	4.33%	0.518%
43	19.377	2738	2744	2751	rBV	4333611	5835461	99.68%	11.900%
44	19.618	2782	2785	2788	rVB	128203	135881	2.32%	0.277%
45	19.736	2795	2805	2813	rBV	3532494	5854359	100.00%	11.939%
46	19.807	2814	2817	2823	rVB2	153464	239328	4.09%	0.488%
47	19.936	2832	2839	2843	rBV2	1125728	1540323	26.31%	3.141%
48	20.230	2886	2889	2892	rVB	213264	220525	3.77%	0.450%
49	20.977	3013	3016	3021	rVB	243658	295952	5.06%	0.604%
50	21.177	3048	3050	3053	rBV	225775	266652	4.55%	0.544%
51	21.483	3097	3102	3107	rBV2	2225551	4618131	78.88%	9.418%
52	21.530	3107	3110	3121	rVB	1529193	2049095	35.00%	4.179%
53	23.177	3385	3390	3395	rBV	1152118	2532638	43.26%	5.165%
54	23.700	3475	3479	3489	rVB	803754	1604481	27.41%	3.272%
55	23.806	3491	3497	3505	rBV	1104748	2332166	39.84%	4.756%
56	23.912	3510	3515	3521	rVB2	954965	1800479	30.75%	3.672%

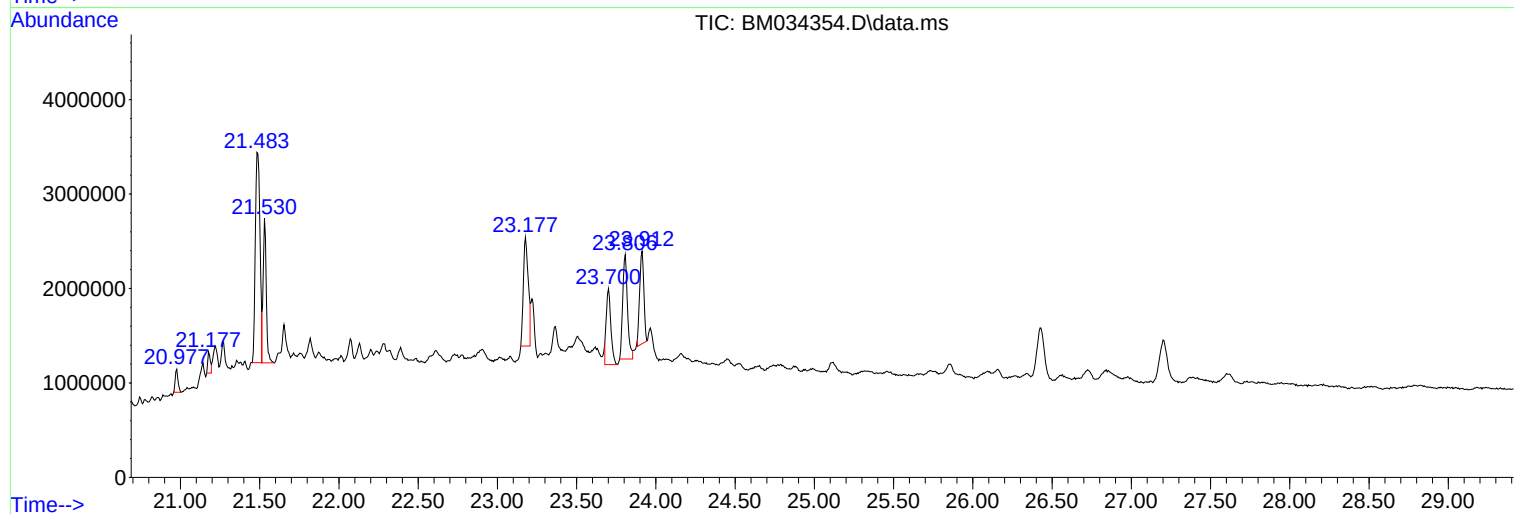
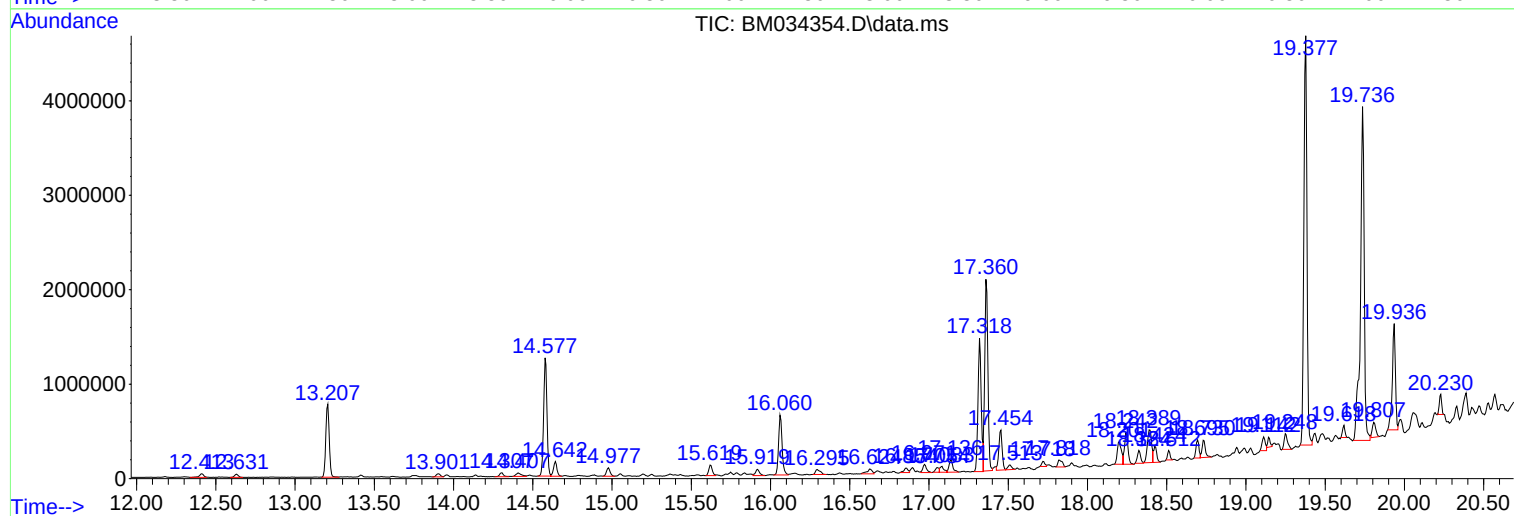
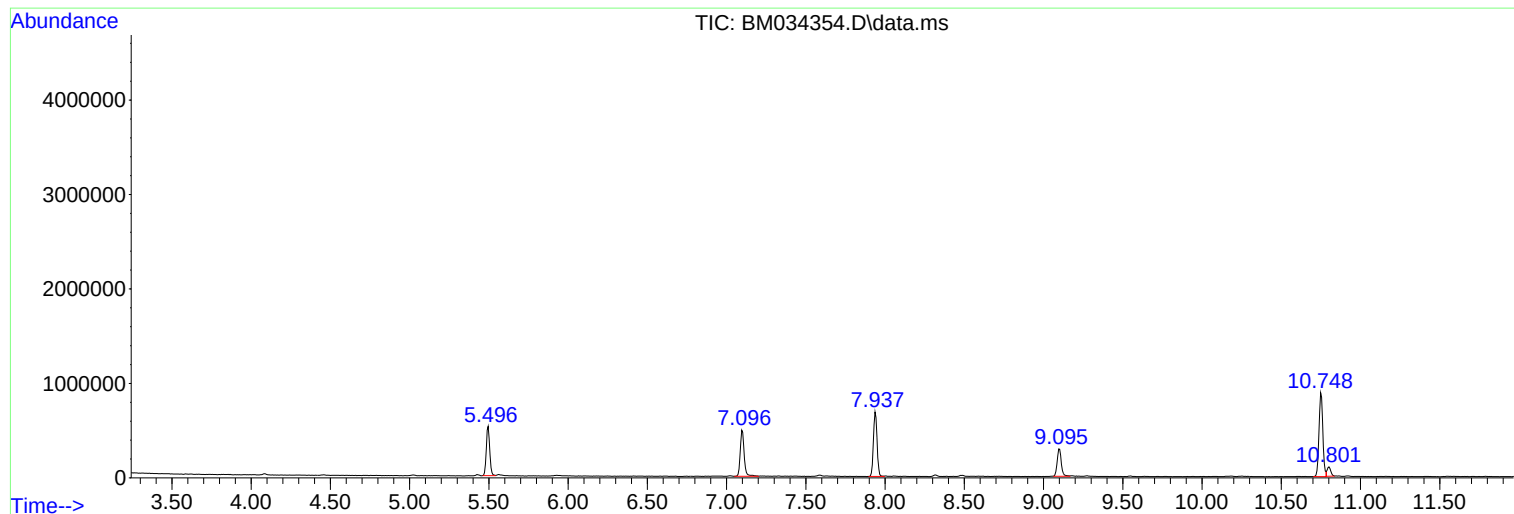
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM031522.M
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P



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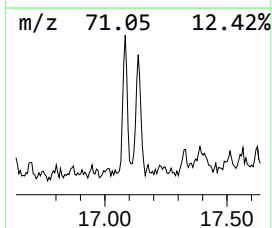
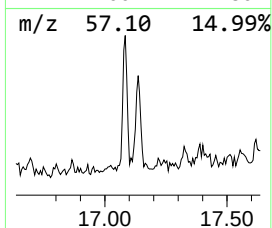
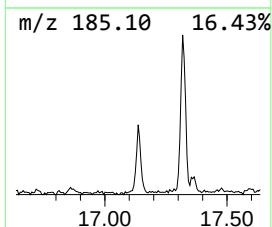
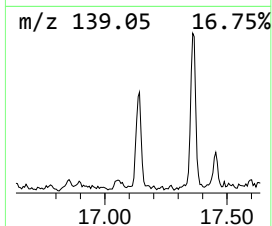
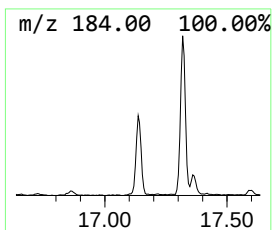
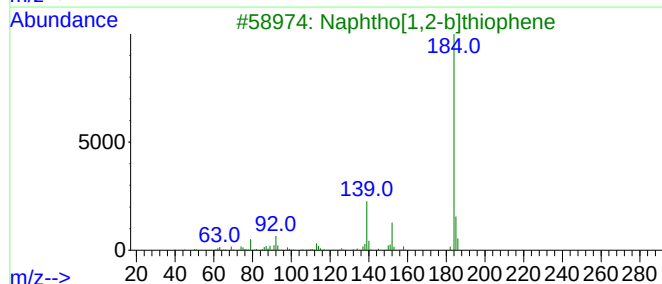
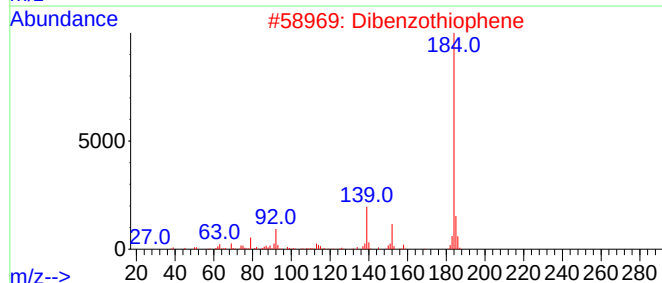
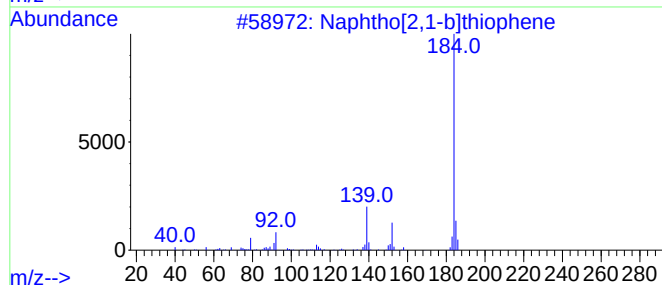
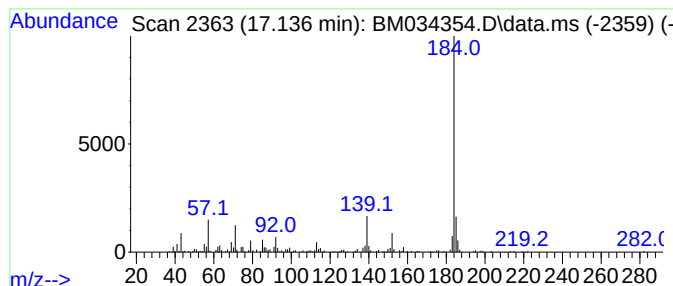
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Naphtho[2,1-b]thiophene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.136	2.10 ng	216741	Phenanthrene-d10	17.318

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphtho[2,1-b]thiophene	184	C12H8S	000233-02-3	95
2			Dibenzothiophene	184	C12H8S	000132-65-0	94
3			Naphtho[1,2-b]thiophene	184	C12H8S	000234-41-3	93
4			Azuleno(2,1-b)thiophene	184	C12H8S	000248-13-5	90
5			Naphtho[2,3-b]thiophene	184	C12H8S	000268-77-9	90



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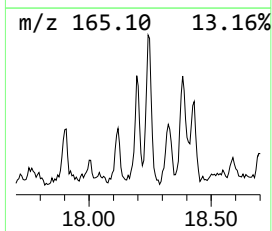
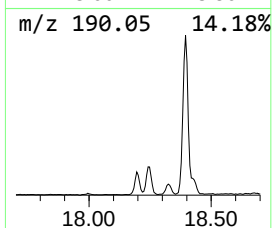
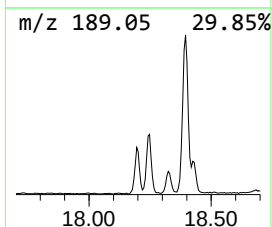
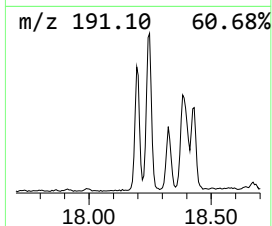
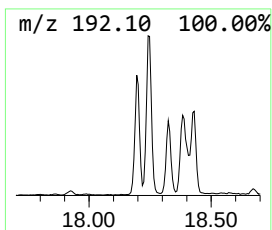
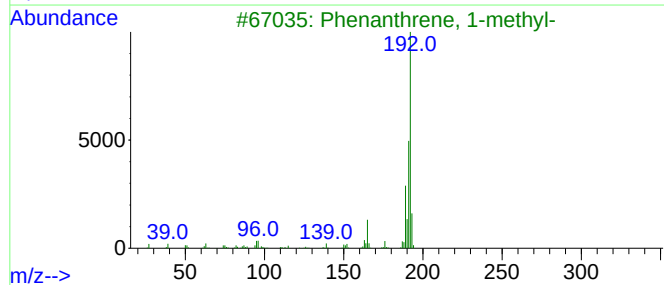
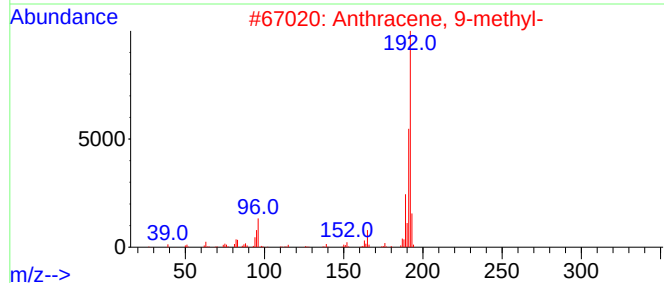
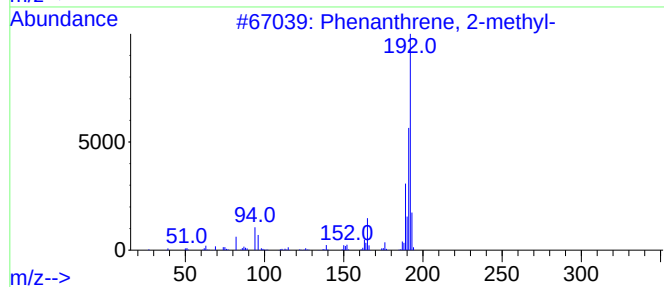
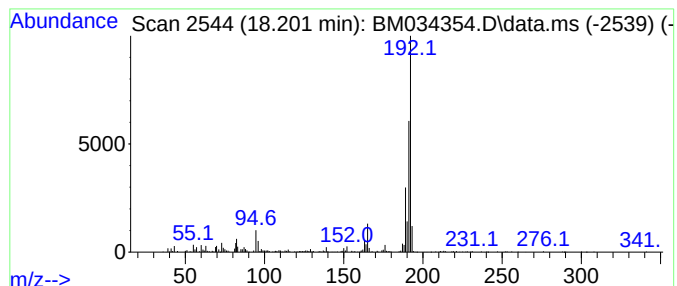
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Phenanthrene, 2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.201	4.32 ng	446232	Phenanthrene-d10	17.318

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	93
2			Anthracene, 9-methyl-	192	C15H12	000779-02-2	91
3			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	90
4			Anthracene, 1-methyl-	192	C15H12	000610-48-0	90
5			Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	90



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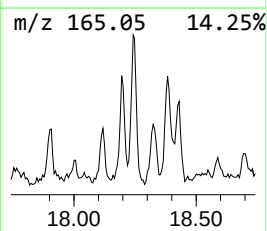
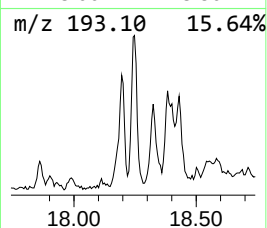
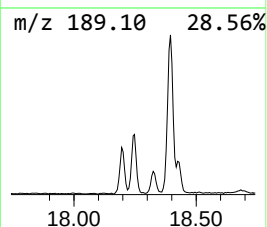
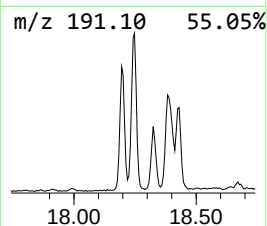
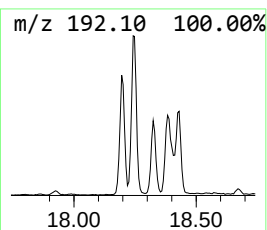
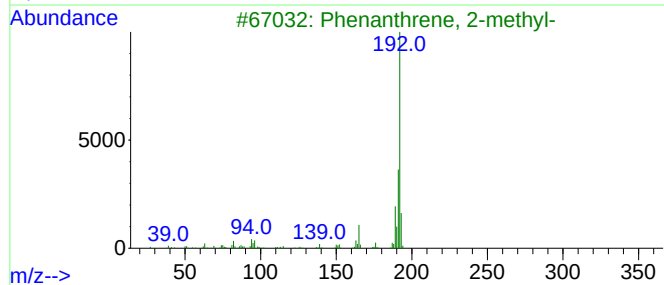
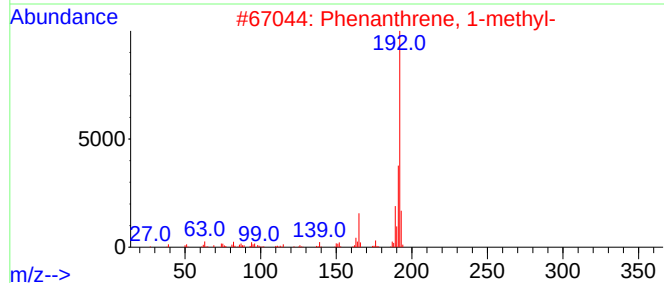
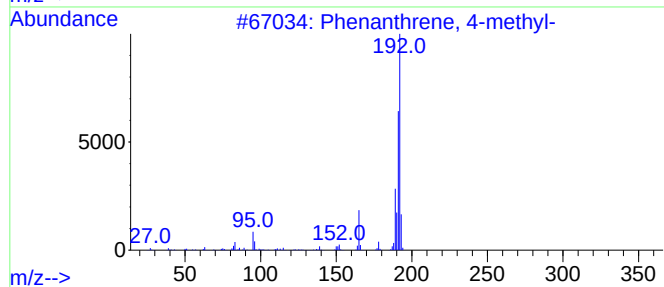
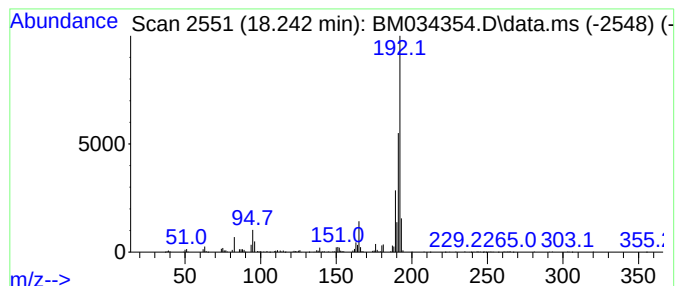
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Peak Number 4 Phenanthrene, 4-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.242	4.85 ng	500815	Phenanthrene-d10	17.318

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	96
2			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	93
3			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	93
4			Anthracene, 1-methyl-	192	C15H12	000610-48-0	93
5			1H-Cyclopropa[1]phenanthrene,1a,...	192	C15H12	000949-41-7	93



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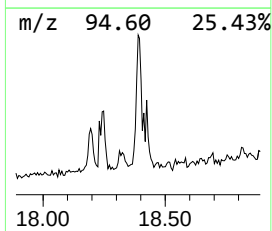
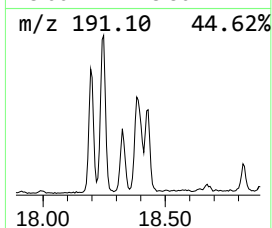
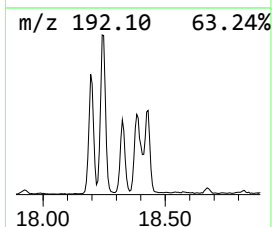
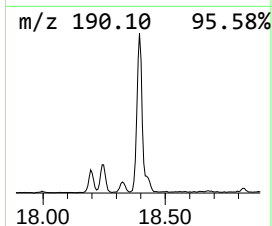
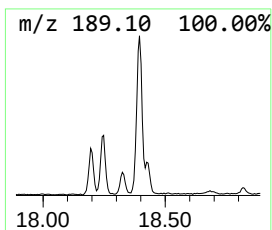
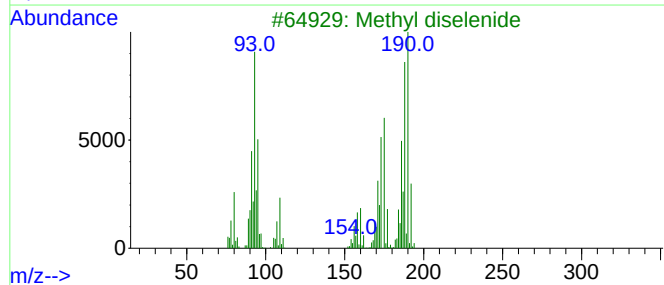
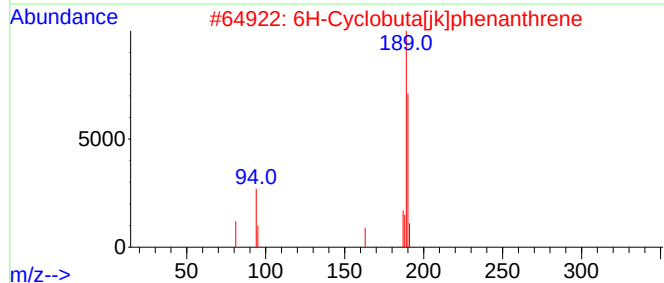
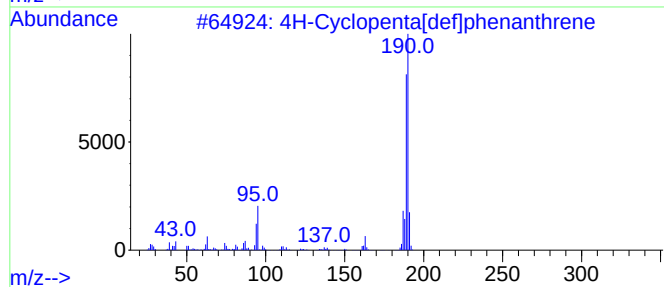
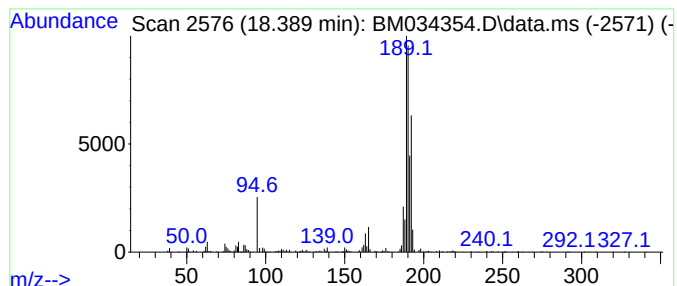
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 4H-Cyclopenta[def]phenanthrene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.389	6.31 ng	651321	Phenanthrene-d10	17.318

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	60
2			6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	58
3			Methyl diselenide	190	C2H6Se2	007101-31-7	41
4			Anthracene, 2-methyl-	192	C15H12	000613-12-7	38
5			5H-Dibenzo[a,d]cycloheptene	192	C15H12	000256-81-5	38



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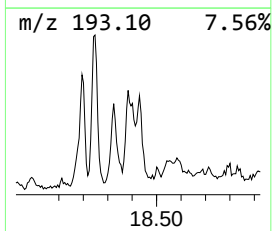
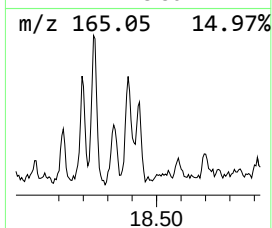
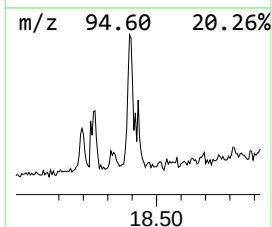
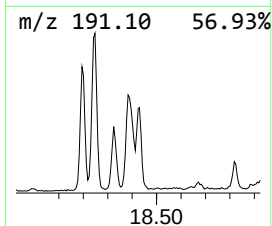
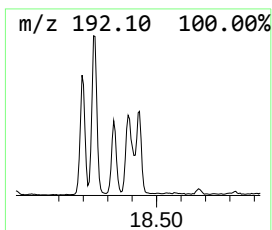
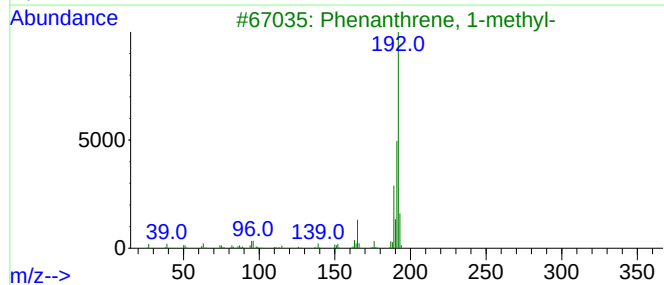
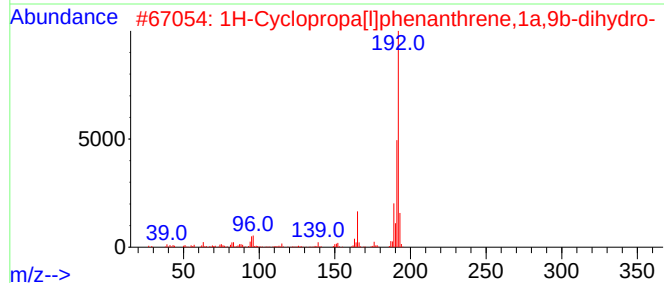
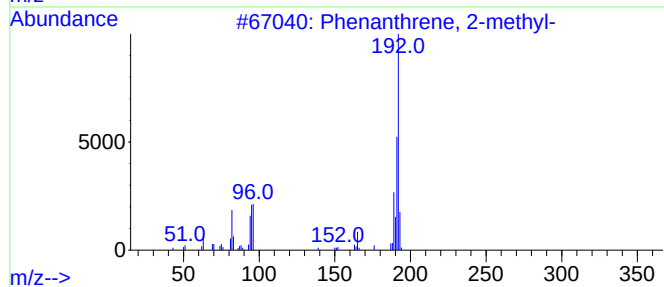
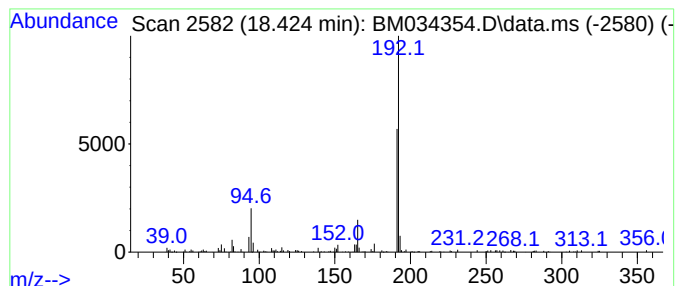
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 1H-Cyclopropa[l]phenanthren... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.424	2.14 ng	220735	Phenanthrene-d10	17.318

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	87
2		1H-Cyclopropa[l]phenanthrene,1a,...	192	C15H12	000949-41-7	87
3		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	87
4		Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	83
5		Anthracene, 9-methyl-	192	C15H12	000779-02-2	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM032322\
Data File : BM034354.D
Acq On : 24 Mar 2022 00:21
Operator : CG/JU
Sample : N1958-08 5X
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C-8-7-8

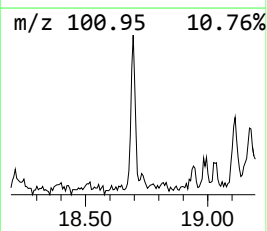
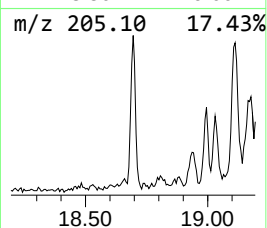
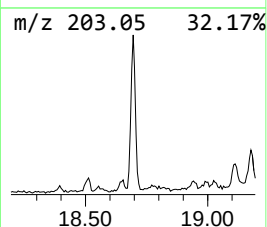
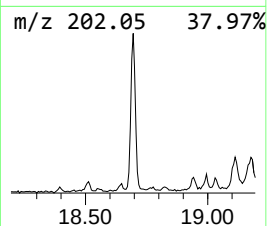
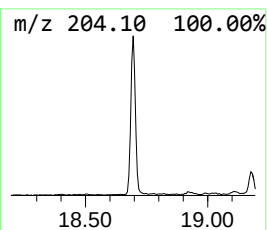
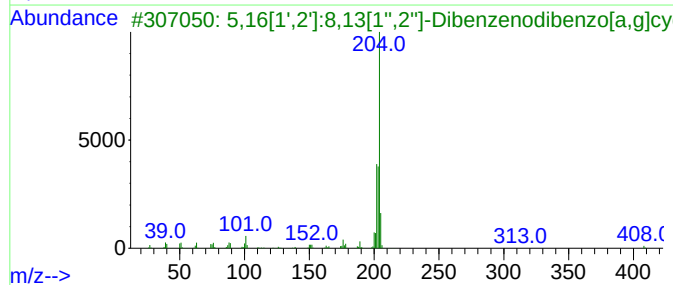
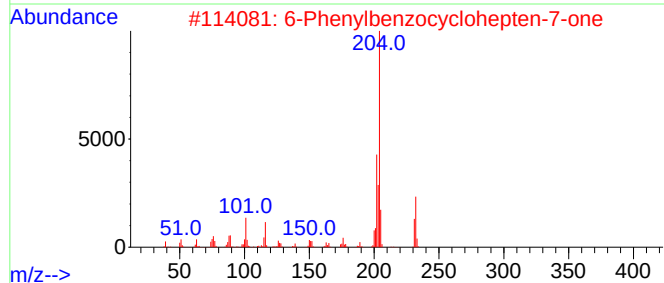
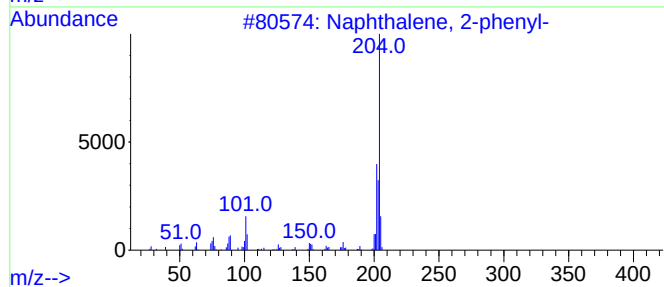
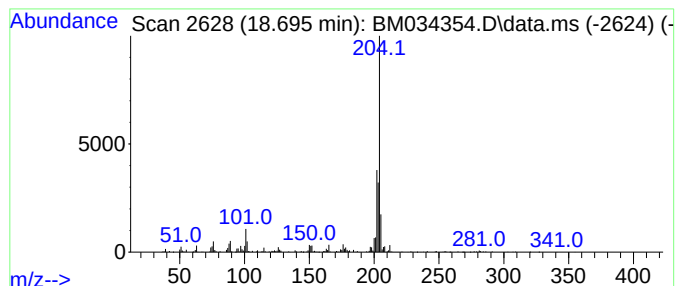
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Naphthalene, 2-phenyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.695	2.42 ng	249597	Phenanthrene-d10	17.318

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	93
2			6-Phenylbenzocyclohepten-7-one	232	C17H12O	093327-56-1	91
3			5,16[1',2']:8,13[1'',2'']-Dibenz...	408	C32H24	005672-97-9	91
4			9,10-Bis(bromomethyl)anthracene	362	C16H12Br2	034373-96-1	70
5			Tricyclo[8.2.2.2(4,7)]hexadeca-2...	204	C16H12	006572-60-7	68



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM032322\
Data File : BM034354.D
Acq On : 24 Mar 2022 00:21
Operator : CG/JU
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Instrument :
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ClientSampleId :
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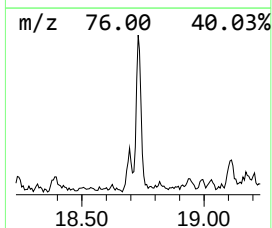
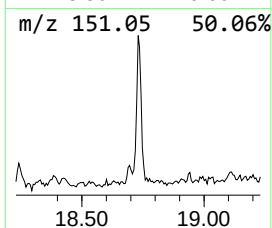
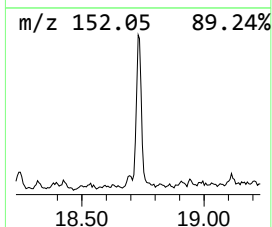
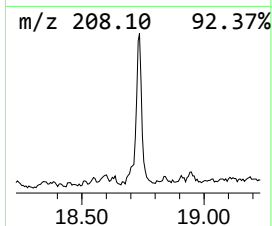
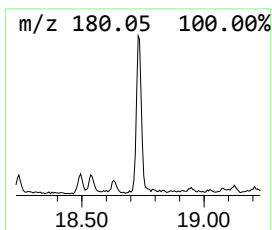
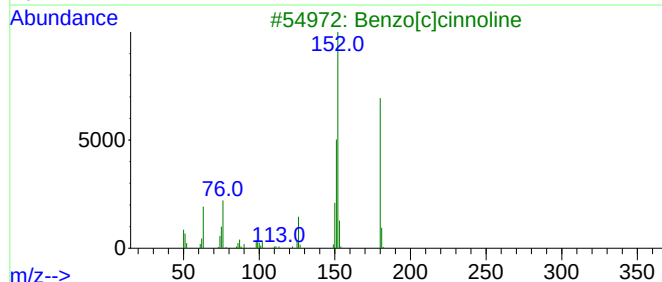
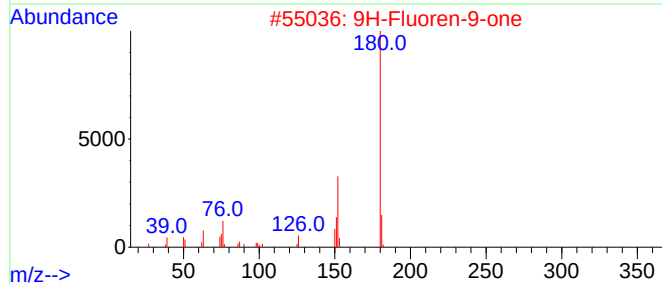
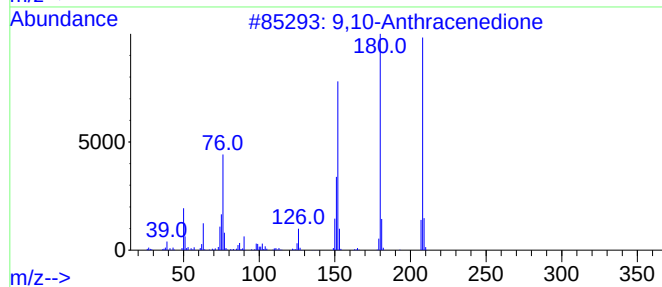
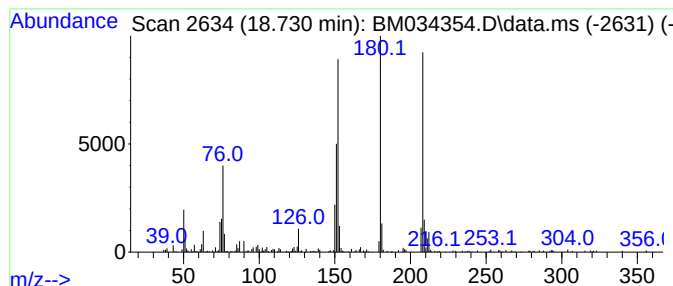
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 9,10-Anthracenedione Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.730	2.50 ng	257909	Phenanthrene-d10	17.318

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		9,10-Anthracenedione	208	C14H8O2	000084-65-1	99
2		9H-Fluoren-9-one	180	C13H8O	000486-25-9	70
3		Benzo[c]cinnoline	180	C12H8N2	000230-17-1	60
4		1H-Phenalen-1-one	180	C13H8O	000548-39-0	60
5		1,4-Anthracenedione	208	C14H8O2	000635-12-1	49



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM032322\
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Sample : N1958-08 5X
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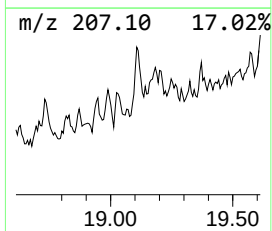
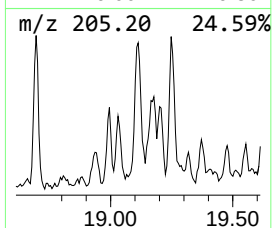
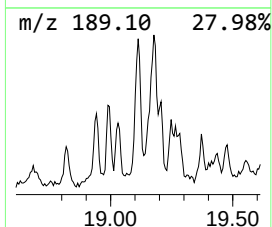
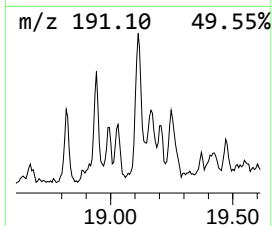
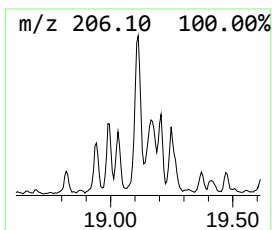
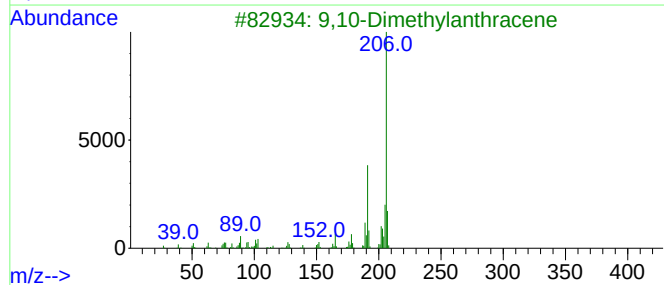
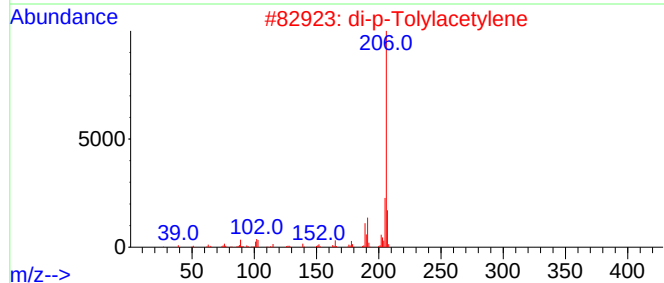
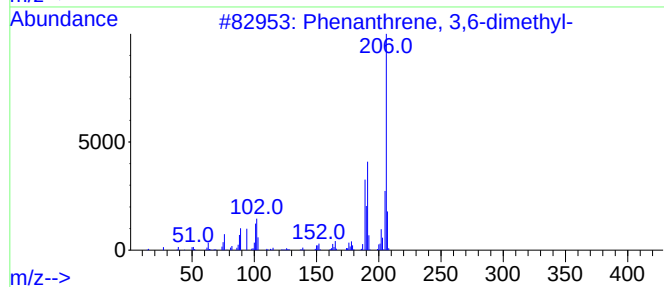
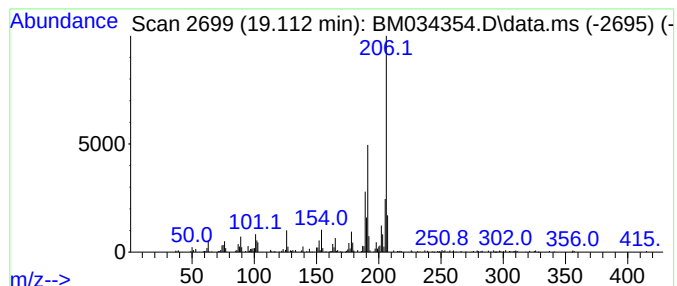
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 Phenanthrene, 3,6-dimethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
19.113	2.07 ng	213330	Phenanthrene-d10	17.318	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 3,6-dimethyl-	206	C16H14	001576-67-6	93
2	di-p-Tolylacetylene	206	C16H14	002789-88-0	91
3	9,10-Dimethylantracene	206	C16H14	000781-43-1	91
4	Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	91
5	Phenanthrene, 1,7-dimethyl-	206	C16H14	000483-87-4	89



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM032322\
Data File : BM034354.D
Acq On : 24 Mar 2022 00:21
Operator : CG/JU
Sample : N1958-08 5X
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C-8-7-8

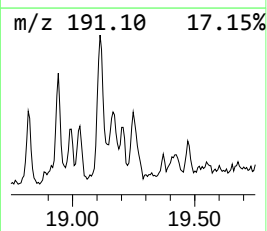
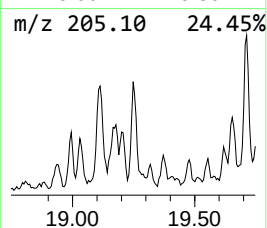
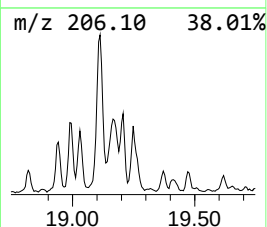
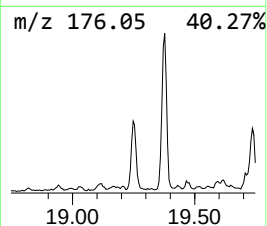
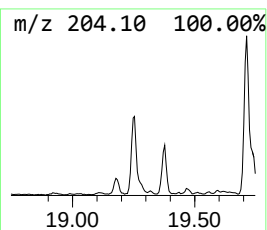
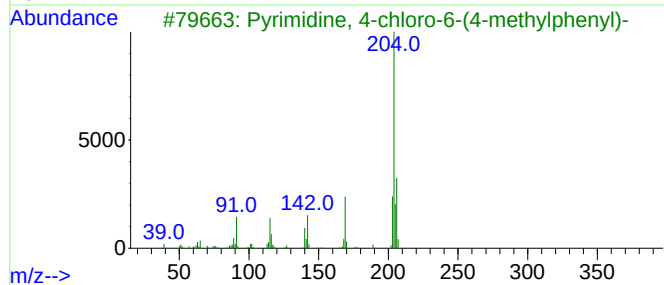
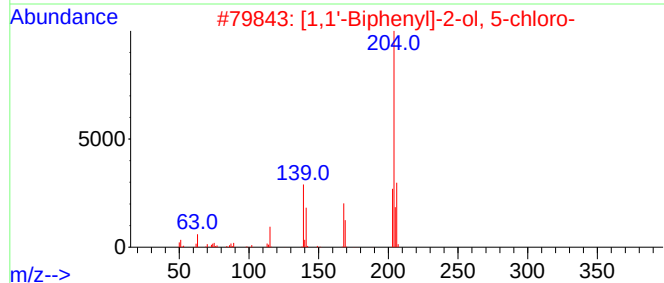
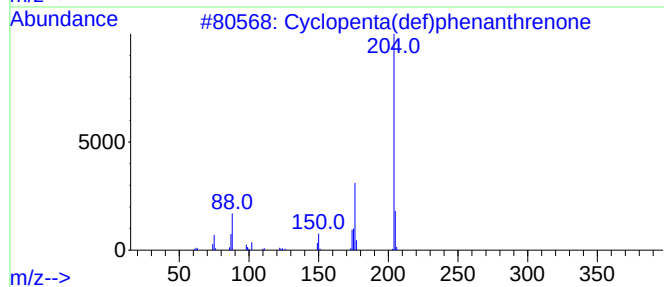
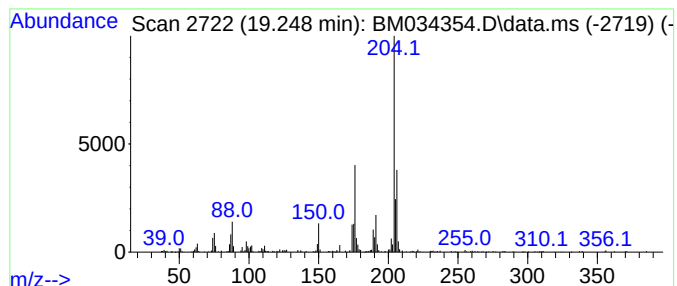
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 Cyclopenta(def)phenanthrenone Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.248	2.46 ng	253766	Phenanthrene-d10	17.318

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopenta(def)phenanthrenone	204	C15H8O	005737-13-3	90
2			[1,1'-Biphenyl]-2-ol, 5-chloro-	204	C12H9ClO	000607-12-5	43
3			Pyrimidine, 4-chloro-6-(4-methyl...	204	C11H9ClN2	1000380-55-8	43
4			1,2,4,8-Tetramethylbicyclo[6.3.0...	204	C15H24	137235-51-9	43
5			Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM032322\
Data File : BM034354.D
Acq On : 24 Mar 2022 00:21
Operator : CG/JU
Sample : N1958-08 5X
Misc :
ALS Vial : 16 Sample Multiplier: 1

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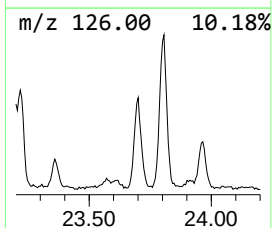
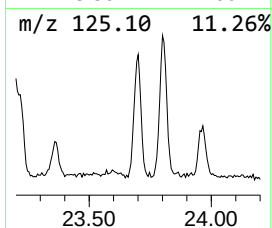
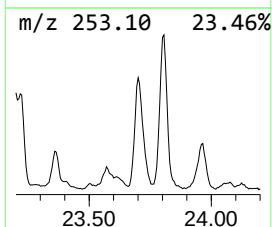
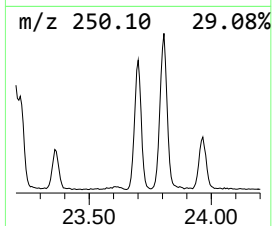
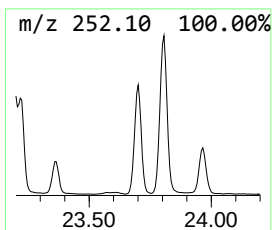
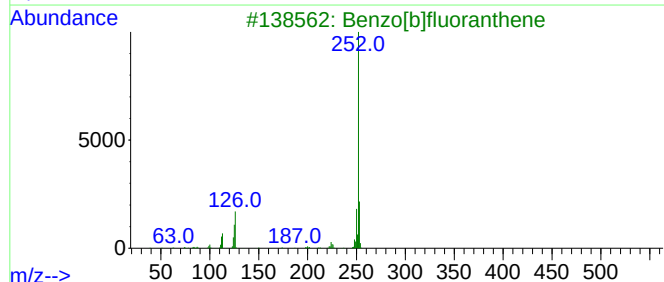
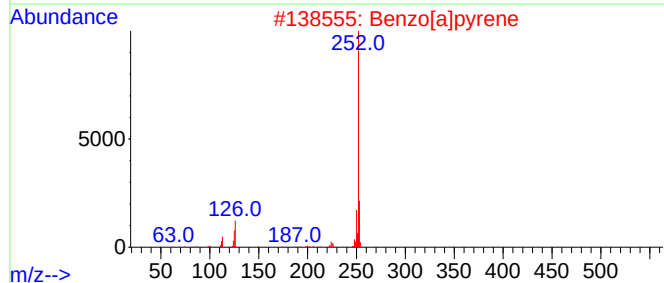
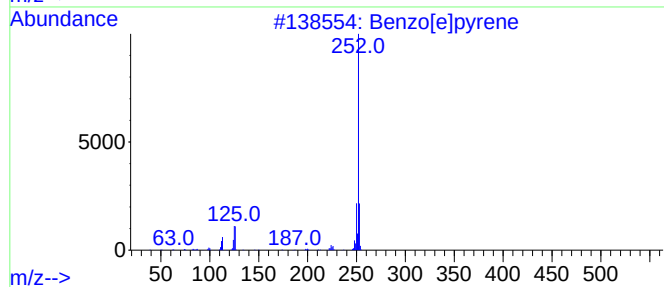
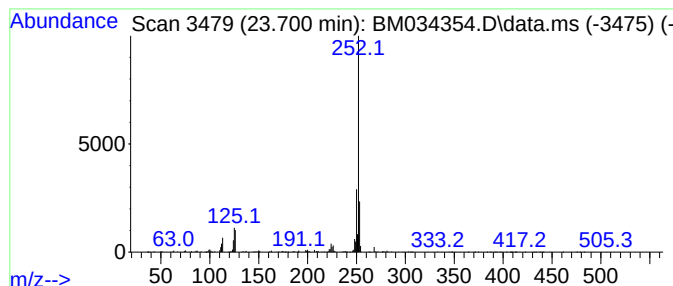
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 Benzo[e]pyrene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.700	17.82 ng	1604480	Perylene-d12	23.912

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[e]pyrene	252	C20H12	000192-97-2	99
2			Benzo[a]pyrene	252	C20H12	000050-32-8	96
3			Benzo[b]fluoranthene	252	C20H12	000205-99-2	95
4			Benzo[k]fluoranthene	252	C20H12	000207-08-9	95
5			Benzo[j]fluoranthene	252	C20H12	000205-82-3	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM032322\
Data File : BM034354.D
Acq On : 24 Mar 2022 00:21
Operator : CG/JU
Sample : N1958-08 5X
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM031522.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
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Phenanthrene, 2...	18.201	4.3	ng	446232	4	17.318	2066000	20.0
Phenanthrene, 4...	18.242	4.8	ng	500815	4	17.318	2066000	20.0
4H-Cyclopenta[d...	18.389	6.3	ng	651321	4	17.318	2066000	20.0
1H-Cyclopropa[1...	18.424	2.1	ng	220735	4	17.318	2066000	20.0
Naphthalene, 2-...	18.695	2.4	ng	249597	4	17.318	2066000	20.0
9,10-Anthracene...	18.730	2.5	ng	257909	4	17.318	2066000	20.0
Phenanthrene, 3...	19.113	2.1	ng	213330	4	17.318	2066000	20.0
Cyclopenta(def)...	19.248	2.5	ng	253766	4	17.318	2066000	20.0
Benzo[e]pyrene	23.700	17.8	ng	1604480	6	23.912	1800480	20.0