

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF092123\
 Data File : BF135380.D
 Acq On : 21 Sep 2023 15:04
 Operator : CG\JU
 Sample : 04501-04
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-3

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_F\Methods\8270-BF091123.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF135380.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.393	558	562	585	rBV	3017719	3235344	100.00%	9.213%
2	6.404	729	734	737	rBV	3108827	3175820	98.16%	9.043%
3	6.757	790	794	801	rBV	1089920	873073	26.99%	2.486%
4	7.322	885	890	893	rBV	2220978	2065347	63.84%	5.881%
5	8.034	1007	1011	1014	rBV	1320659	1084141	33.51%	3.087%
6	8.845	1145	1149	1152	rBV	50554	41332	1.28%	0.118%
7	9.116	1190	1195	1198	rBV	3445346	3101510	95.86%	8.832%
8	9.381	1236	1240	1243	rBV2	27686	37532	1.16%	0.107%
9	9.451	1248	1252	1254	rBV	70809	66033	2.04%	0.188%
10	9.475	1254	1256	1261	rVB2	29832	37681	1.16%	0.107%
11	9.569	1269	1272	1278	rVB	36598	43884	1.36%	0.125%
12	9.651	1283	1286	1291	rBV	174787	144424	4.46%	0.411%
13	9.792	1303	1310	1313	rBV	1217288	1053036	32.55%	2.999%
14	9.822	1313	1315	1319	rVB	71718	63017	1.95%	0.179%
15	9.998	1340	1345	1348	rBV2	49343	53769	1.66%	0.153%
16	10.034	1348	1351	1355	rVB	46041	37348	1.15%	0.106%
17	10.110	1360	1364	1367	rBV2	52254	57567	1.78%	0.164%
18	10.198	1375	1379	1381	rBV2	35838	45681	1.41%	0.130%
19	10.339	1400	1403	1409	rVB2	89480	102730	3.18%	0.293%
20	10.586	1441	1445	1457	rVB	1565193	1552266	47.98%	4.420%
21	10.710	1462	1466	1470	rVB3	59589	74074	2.29%	0.211%
22	10.892	1492	1497	1499	rBV3	50562	62930	1.95%	0.179%
23	10.922	1499	1502	1504	rVB2	39887	37112	1.15%	0.106%
24	11.092	1528	1531	1534	rVB2	55869	48371	1.50%	0.138%
25	11.181	1543	1546	1552	rVB	97265	93923	2.90%	0.267%
26	11.286	1560	1564	1566	rBV	947018	931629	28.80%	2.653%
27	11.310	1566	1568	1571	rVV	926861	746809	23.08%	2.127%
28	11.363	1574	1577	1579	rVB	137358	120551	3.73%	0.343%
29	11.398	1579	1583	1585	rBV2	44372	45657	1.41%	0.130%
30	11.522	1600	1604	1609	rBV2	63251	91945	2.84%	0.262%
31	11.580	1612	1614	1616	rVV	74538	57378	1.77%	0.163%
32	11.616	1617	1620	1625	rVB2	97059	103276	3.19%	0.294%
33	11.698	1628	1634	1638	rVB2	61141	96209	2.97%	0.274%
34	11.745	1639	1642	1645	rBV	38441	41769	1.29%	0.119%
35	11.786	1646	1649	1652	rVV	318067	303831	9.39%	0.865%
36	11.816	1652	1654	1659	rVV2	532641	544796	16.84%	1.551%

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37	11.863	1659	1662	1665	rVV2	73904	93275	2.88%	0.266%
38	11.898	1665	1668	1670	rVV2	408913	398075	12.30%	1.134%
39	11.922	1670	1672	1678	rVB	245121	219991	6.80%	0.626%
40	11.992	1678	1684	1687	rBV3	40635	81235	2.51%	0.231%
41	12.022	1687	1689	1693	rVV	61805	55267	1.71%	0.157%
42	12.086	1697	1700	1702	rVV2	160326	155147	4.80%	0.442%
43	12.116	1702	1705	1710	rVB	176199	174939	5.41%	0.498%
44	12.233	1723	1725	1728	rVV	83961	77403	2.39%	0.220%
45	12.269	1728	1731	1733	rVV	126429	107688	3.33%	0.307%
46	12.292	1733	1735	1738	rVB	97250	75675	2.34%	0.215%
47	12.339	1738	1743	1746	rBV	309799	328130	10.14%	0.934%
48	12.375	1746	1749	1752	rVV2	151590	195773	6.05%	0.557%
49	12.398	1752	1753	1755	rVB	101860	54687	1.69%	0.156%
50	12.433	1755	1759	1762	rBV3	144298	192581	5.95%	0.548%
51	12.504	1767	1771	1775	rBV	1220035	1032038	31.90%	2.939%
52	12.551	1776	1779	1780	rBV	68731	55443	1.71%	0.158%
53	12.569	1780	1782	1785	rVV2	58040	50852	1.57%	0.145%
54	12.604	1785	1788	1792	rVV2	83971	123825	3.83%	0.353%
55	12.657	1792	1797	1800	rVV3	79638	132950	4.11%	0.379%
56	12.733	1804	1810	1813	rBV	1298421	1274333	39.39%	3.629%
57	12.792	1817	1820	1823	rVB2	129299	135903	4.20%	0.387%
58	12.851	1826	1830	1831	rBV2	46830	50264	1.55%	0.143%
59	12.880	1831	1835	1842	rVB	2440183	2288231	70.73%	6.516%
60	12.951	1843	1847	1855	rBV4	162913	273658	8.46%	0.779%
61	13.010	1855	1857	1859	rBV	51582	38490	1.19%	0.110%
62	13.045	1859	1863	1864	rBV3	133356	153051	4.73%	0.436%
63	13.063	1864	1866	1870	rVB	219459	221396	6.84%	0.630%
64	13.110	1871	1874	1876	rBV3	49558	58095	1.80%	0.165%
65	13.133	1876	1878	1880	rVV	131035	96487	2.98%	0.275%
66	13.169	1880	1884	1888	rVV2	239932	220337	6.81%	0.627%
67	13.263	1897	1900	1902	rVV	204788	153914	4.76%	0.438%
68	13.292	1902	1905	1908	rVV	194561	162772	5.03%	0.463%
69	13.386	1918	1921	1924	rVB3	37372	42463	1.31%	0.121%
70	13.463	1932	1934	1937	rBV3	27805	39788	1.23%	0.113%
71	13.551	1947	1949	1951	rBV3	37167	39570	1.22%	0.113%
72	13.580	1951	1954	1958	rVB	142279	150170	4.64%	0.428%
73	13.639	1961	1964	1967	rVB	63878	50624	1.56%	0.144%
74	13.692	1967	1973	1975	rBV2	174222	255594	7.90%	0.728%
75	13.716	1975	1977	1979	rVV	132412	103140	3.19%	0.294%
76	13.739	1979	1981	1983	rVV	88442	86753	2.68%	0.247%

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77	13.792	1987	1990	1993	rVB	147077	133283	4.12%	0.380%
78	13.863	1999	2002	2003	rBV	41546	39770	1.23%	0.113%
79	13.939	2010	2015	2018	rBV2	963649	1346313	41.61%	3.834%
80	13.969	2018	2020	2028	rVB	603506	538976	16.66%	1.535%
81	14.027	2028	2030	2032	rBV	72833	61966	1.92%	0.176%
82	14.092	2039	2041	2042	rBV	68931	53707	1.66%	0.153%
83	14.198	2057	2059	2063	rVB3	42737	49162	1.52%	0.140%
84	14.263	2067	2070	2073	rBV3	44330	53279	1.65%	0.152%
85	14.298	2073	2076	2077	rBV	60086	58235	1.80%	0.166%
86	14.333	2077	2082	2085	rVB	203937	224507	6.94%	0.639%
87	14.363	2085	2087	2091	rBV	118374	107577	3.33%	0.306%
88	14.457	2100	2103	2105	rBV	131272	118941	3.68%	0.339%
89	14.480	2105	2107	2110	rVB2	76636	64667	2.00%	0.184%
90	14.527	2113	2115	2118	rVB2	54618	48363	1.49%	0.138%
91	14.721	2145	2148	2151	rBV2	55780	73253	2.26%	0.209%
92	14.827	2163	2166	2169	rBV4	49576	53970	1.67%	0.154%
93	14.980	2188	2192	2196	rBV	326227	508071	15.70%	1.447%
94	15.092	2208	2211	2214	rVB	68215	71009	2.19%	0.202%
95	15.280	2240	2243	2249	rVB	224898	283652	8.77%	0.808%
96	15.345	2250	2254	2259	rVV	287713	357479	11.05%	1.018%
97	15.410	2260	2265	2268	rVV	490811	616491	19.05%	1.755%
98	15.439	2268	2270	2277	rVB2	88773	114206	3.53%	0.325%
99	16.886	2511	2516	2523	rVB2	93791	194074	6.00%	0.553%
100	17.333	2587	2592	2603	rVB2	79151	175585	5.43%	0.500%

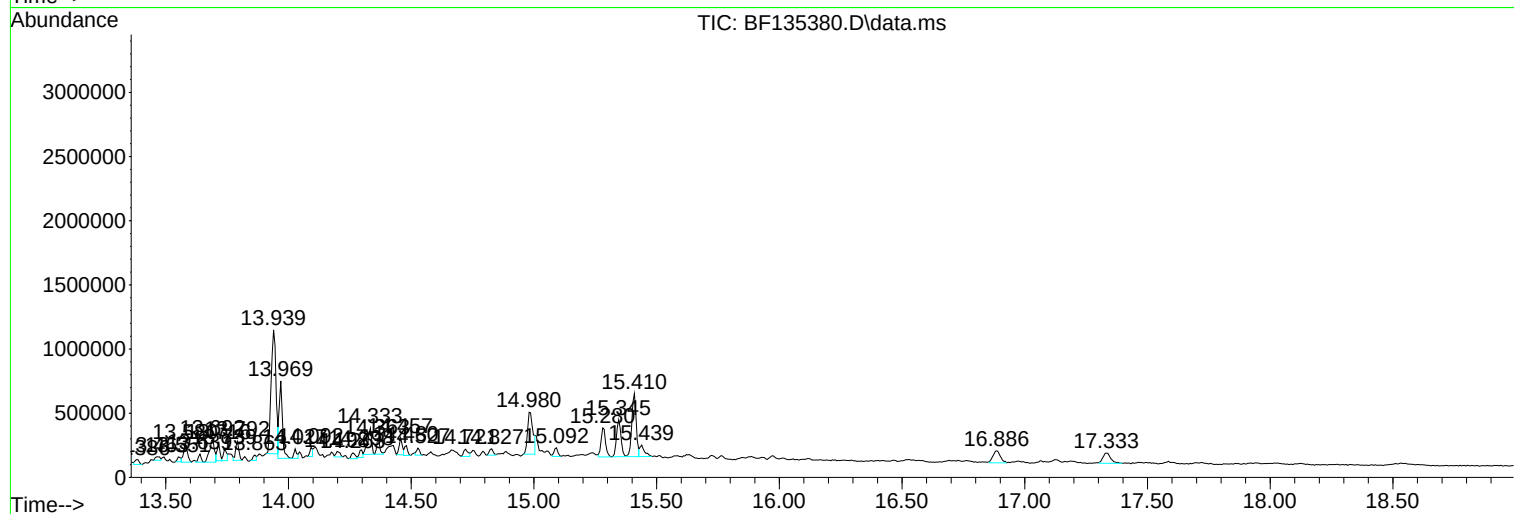
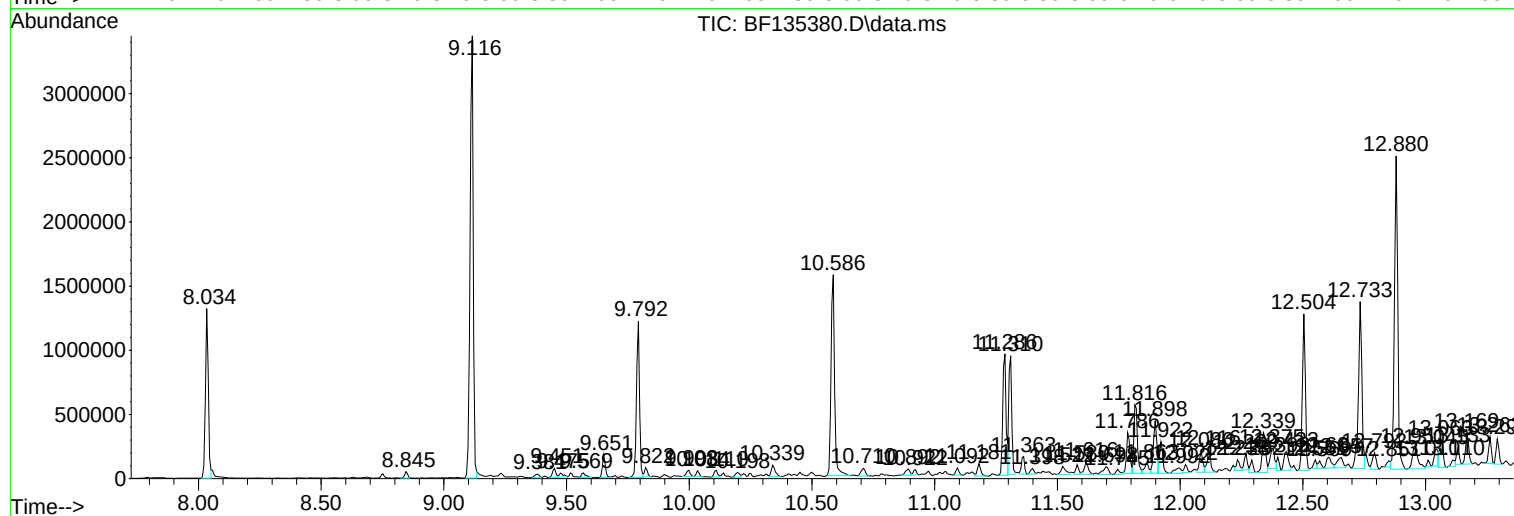
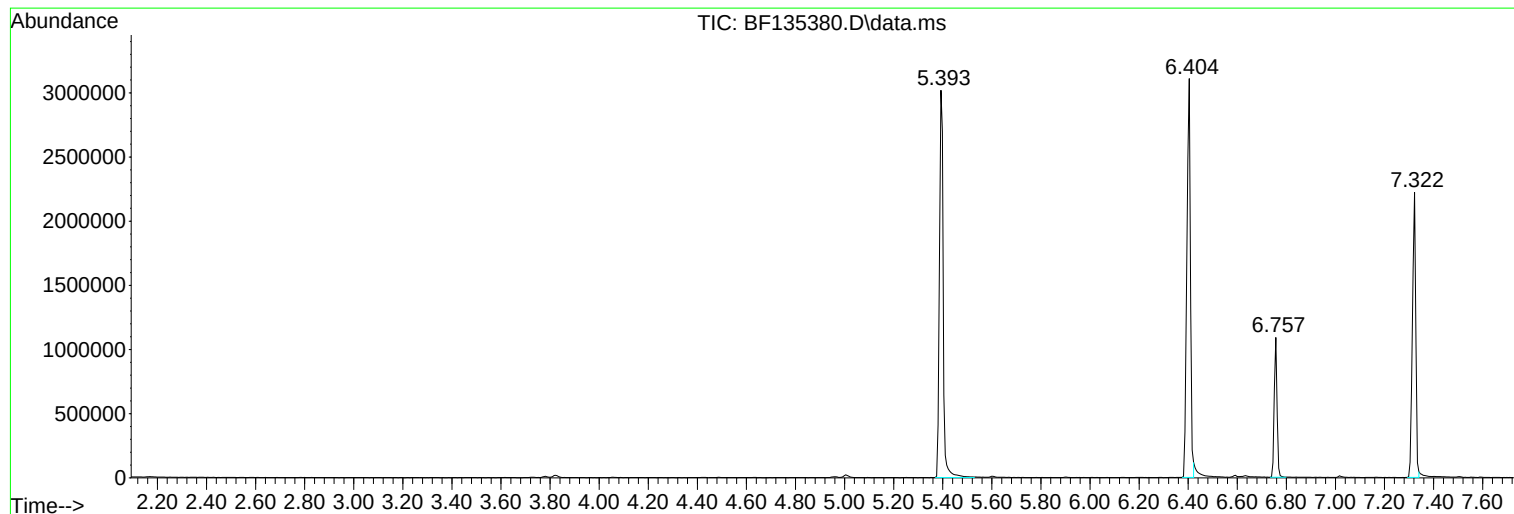
Sum of corrected areas: 35118338

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Instrument :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF091123.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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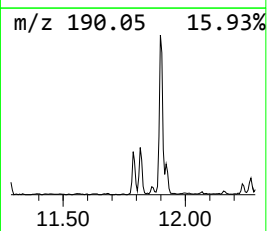
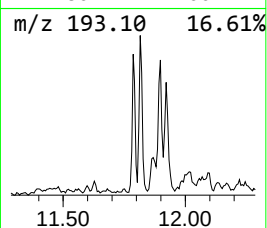
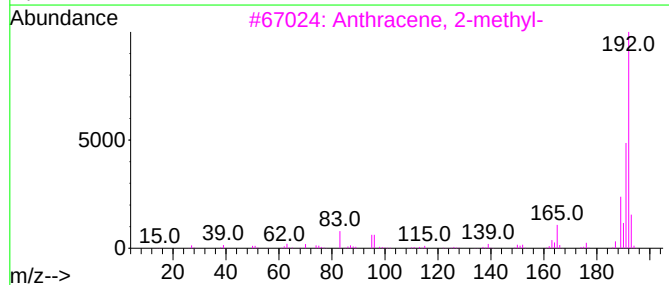
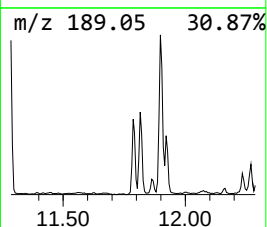
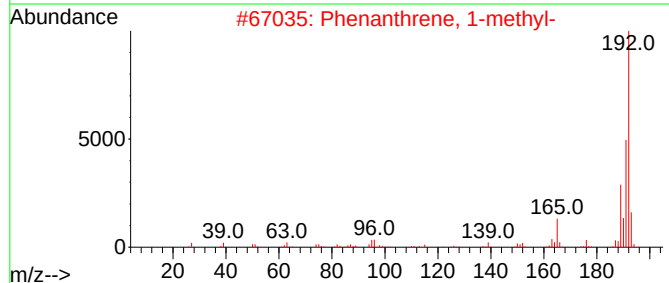
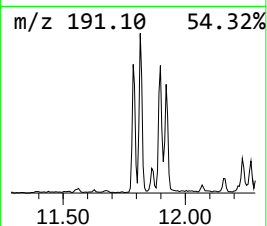
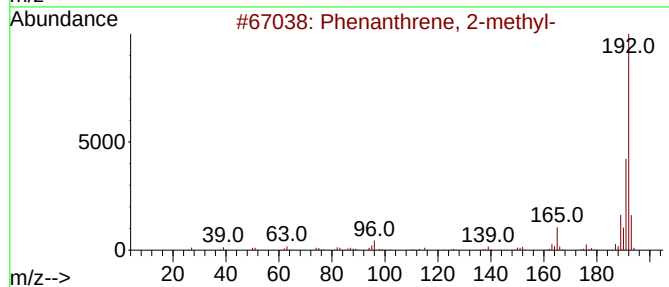
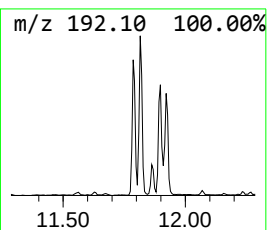
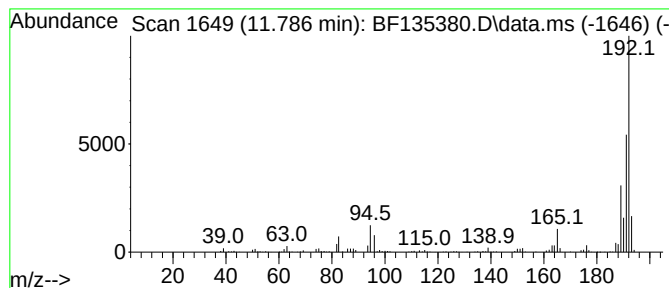
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF091123.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.786	6.52 ng	303831	Phenanthrene-d10	11.286

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	96
2			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	96
3			Anthracene, 2-methyl-	192	C15H12	000613-12-7	95
4			Anthracene, 9-methyl-	192	C15H12	000779-02-2	91
5			Anthracene, 1-methyl-	192	C15H12	000610-48-0	90



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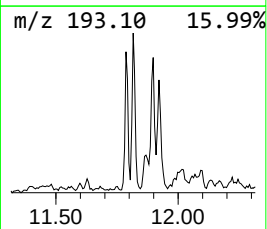
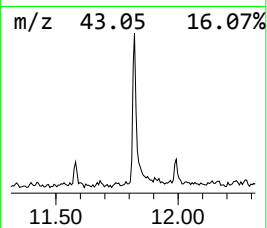
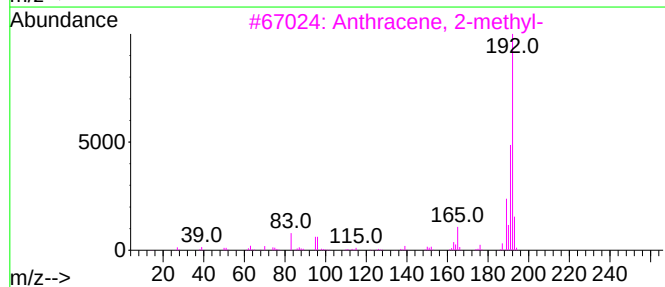
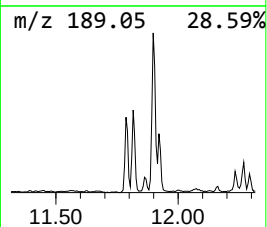
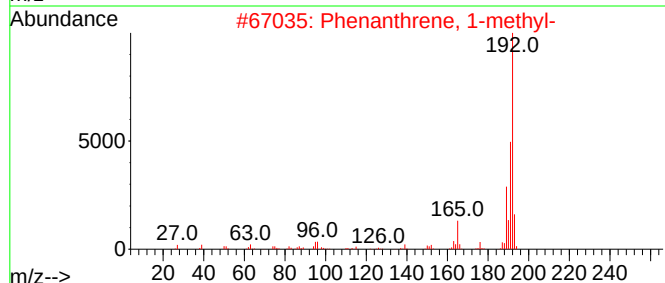
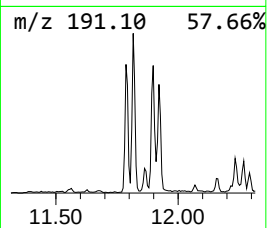
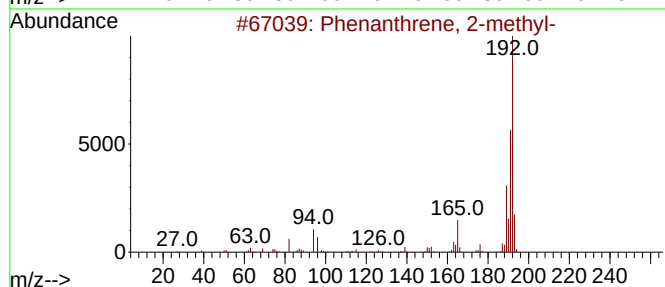
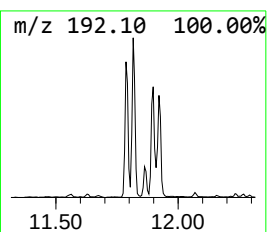
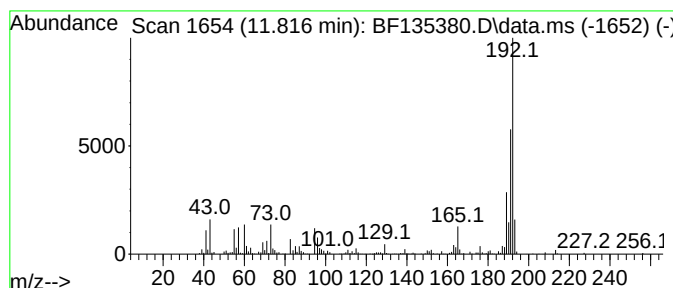
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.816	11.70 ng	544796	Phenanthrene-d10	11.286	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	98
2	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	96
3	Anthracene, 2-methyl-	192	C15H12	000613-12-7	96
4	Anthracene, 9-methyl-	192	C15H12	000779-02-2	95
5	1H-Cyclopropa[1]phenanthrene,1a,...	192	C15H12	000949-41-7	95



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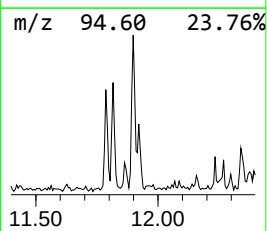
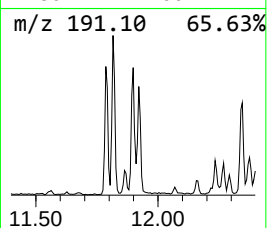
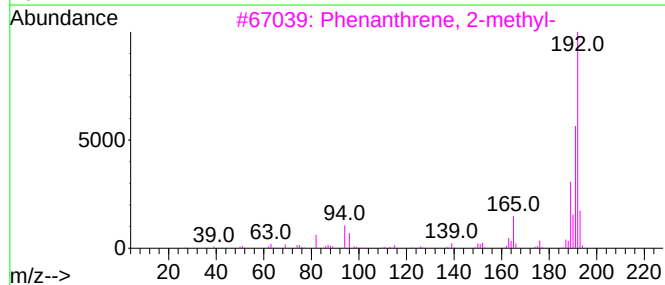
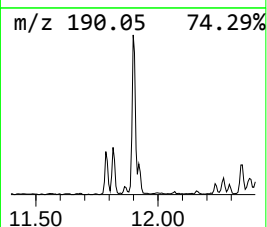
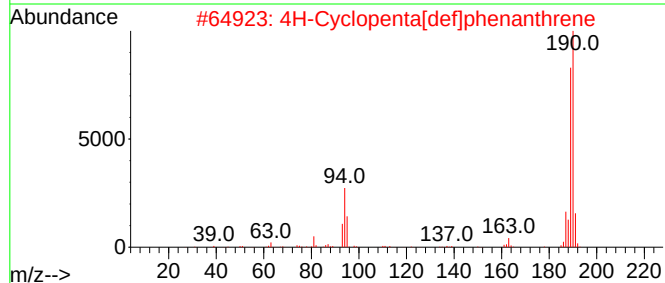
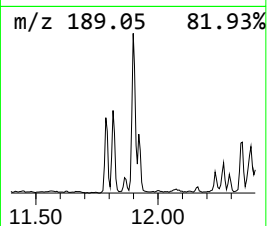
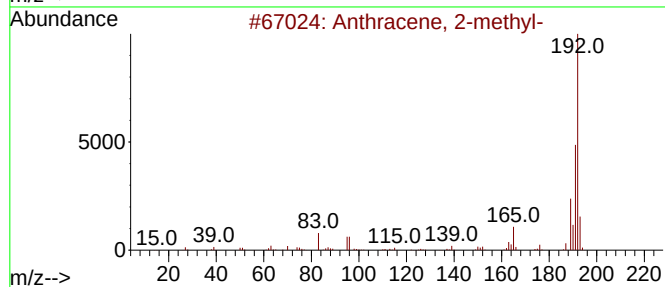
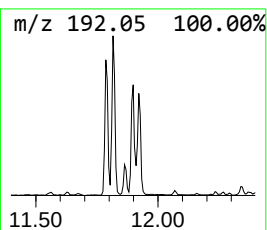
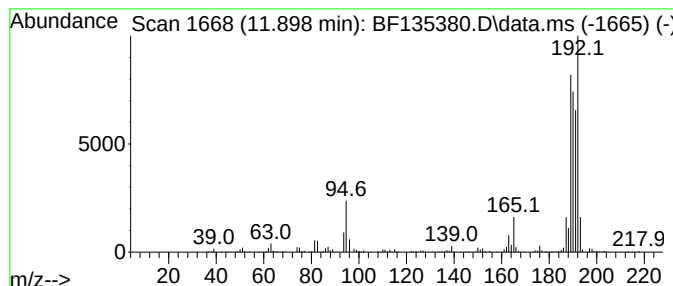
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.898	8.55 ng	398075	Phenanthrene-d10	11.286

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Anthracene, 2-methyl-	192	C15H12	000613-12-7	83
2			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	60
3			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	46
4			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	46
5			Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	42



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF092123\
Data File : BF135380.D
Acq On : 21 Sep 2023 15:04
Operator : CG\JU
Sample : 04501-04
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
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ClientSampleId :
TP-3

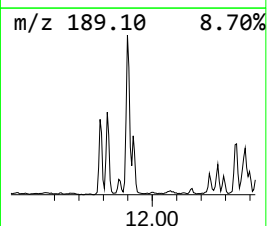
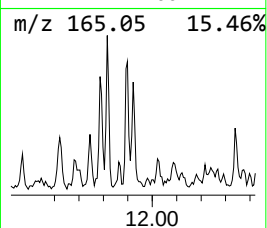
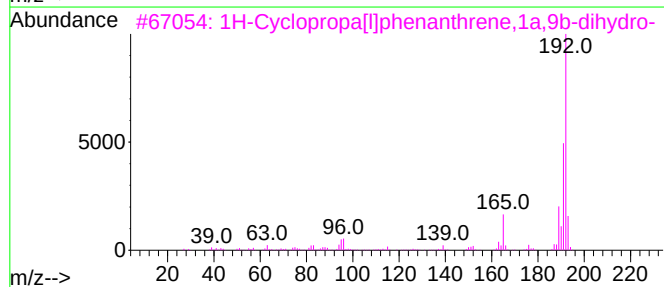
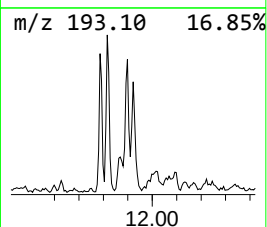
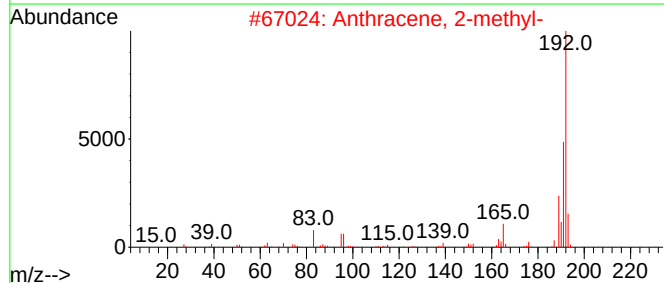
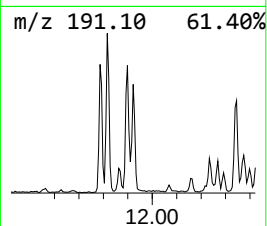
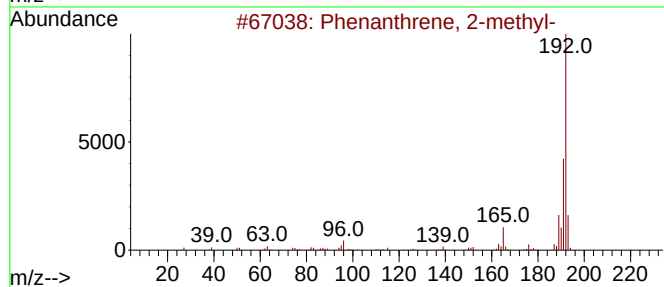
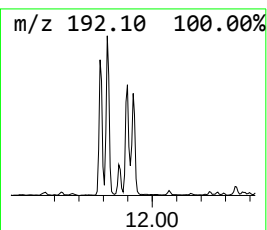
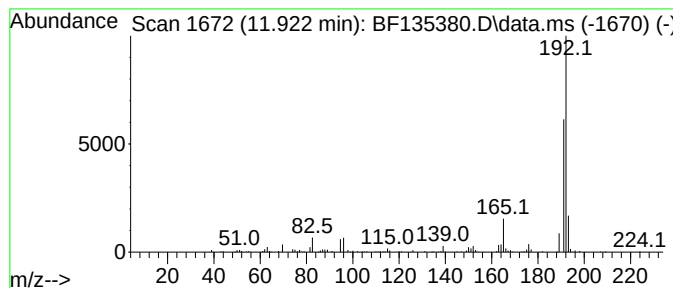
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 4 1H-Cyclopropa[l]phenanthren... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.922	4.72 ng	219991	Phenanthrene-d10	11.286

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	95
2		Anthracene, 2-methyl-	192	C15H12	000613-12-7	94
3		1H-Cyclopropa[l]phenanthrene,1a,...	192	C15H12	000949-41-7	91
4		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	91
5		Anthracene, 9-methyl-	192	C15H12	000779-02-2	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF092123\
Data File : BF135380.D
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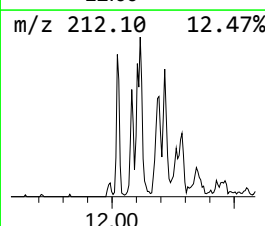
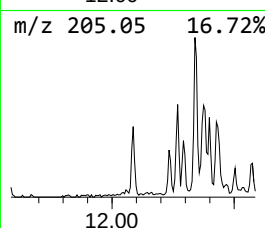
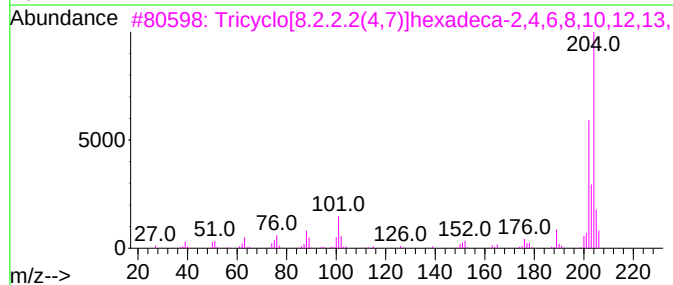
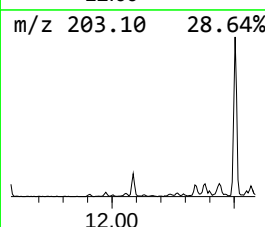
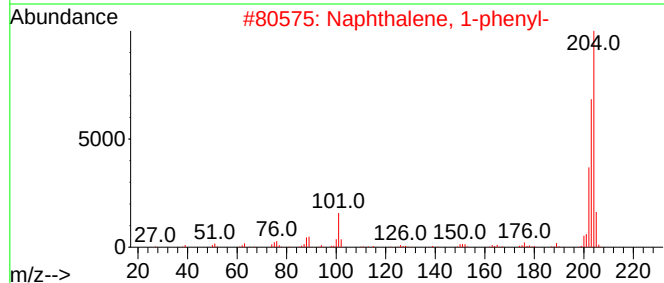
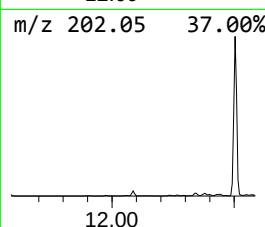
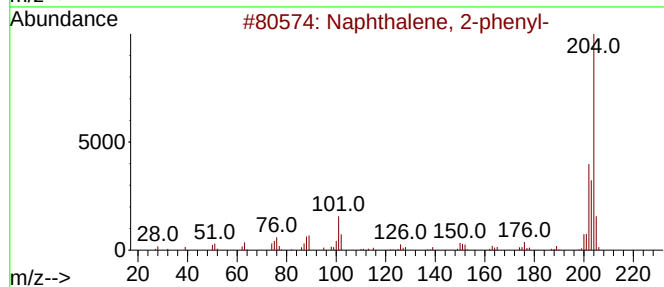
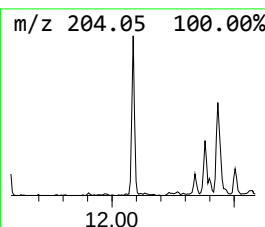
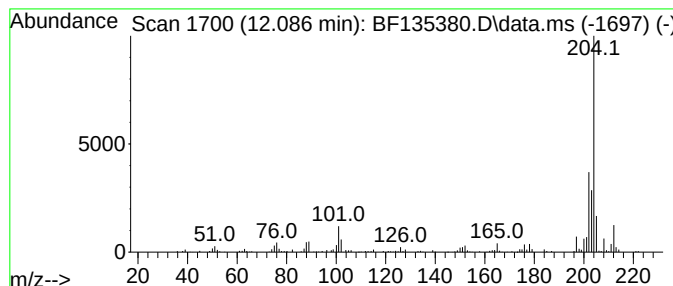
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Naphthalene, 2-phenyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.086	3.33 ng	155147	Phenanthrene-d10	11.286

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	89
2			Naphthalene, 1-phenyl-	204	C16H12	000605-02-7	78
3			Tricyclo[8.2.2.2(4,7)]hexadeca-2...	204	C16H12	006572-60-7	76
4			5,16[1',2'] :8,13[1',2'']-Dibenz...	408	C32H24	005672-97-9	72
5			5H-Dibenzo[a,d]cycloheptene, 5-m...	204	C16H12	002975-79-3	60



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF092123\
Data File : BF135380.D
Acq On : 21 Sep 2023 15:04
Operator : CG\JU
Sample : 04501-04
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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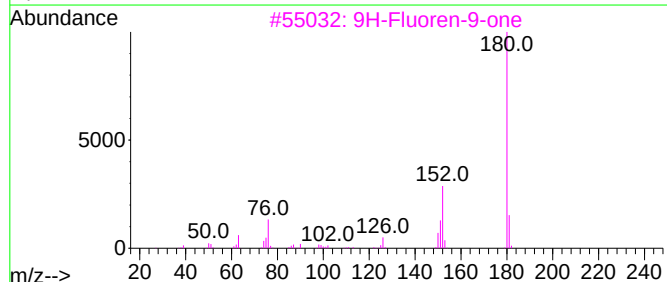
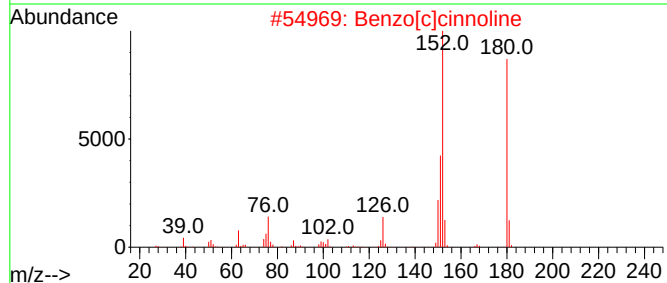
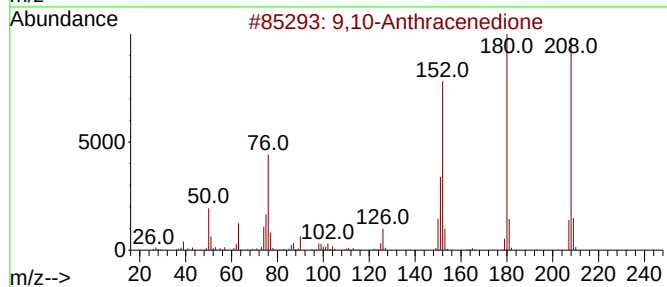
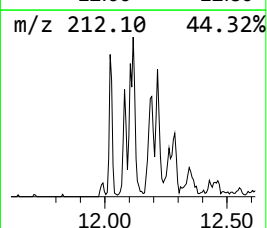
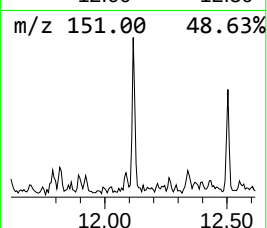
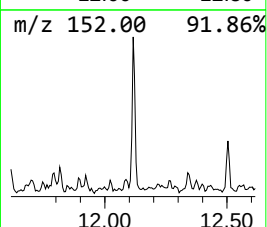
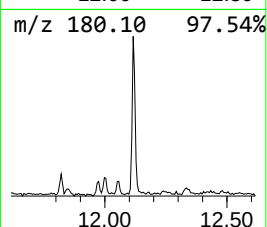
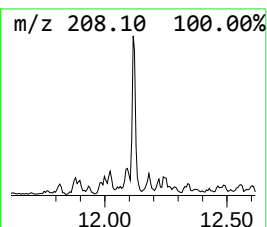
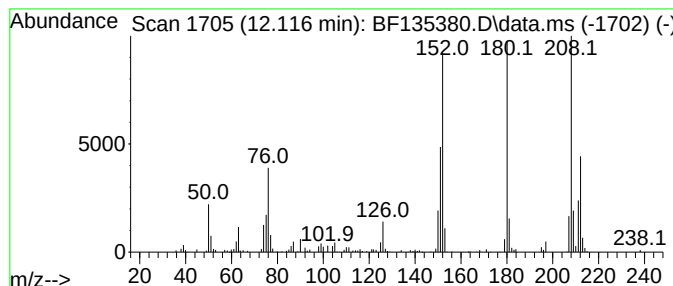
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 9,10-Anthracenedione Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.116	3.76 ng	174939	Phenanthrene-d10	11.286

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		9,10-Anthracenedione	208	C14H8O2	000084-65-1	99
2		Benzo[c]cinnoline	180	C12H8N2	000230-17-1	89
3		9H-Fluoren-9-one	180	C13H8O	000486-25-9	70
4		1H-Phenalen-1-one	180	C13H8O	000548-39-0	50
5		9,10-Phenanthrenedione	208	C14H8O2	000084-11-7	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF092123\
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ClientSampleId :
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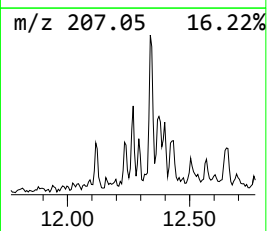
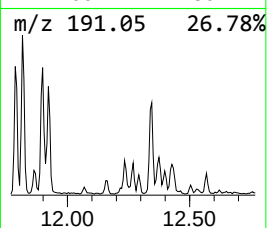
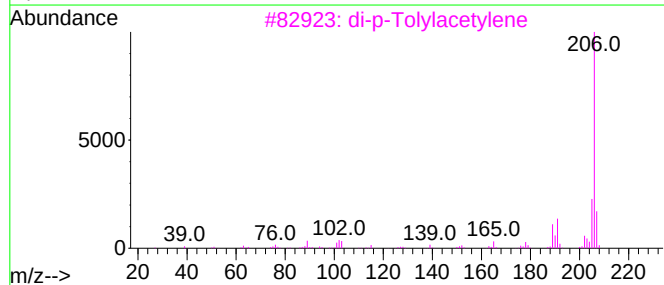
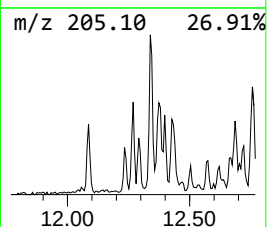
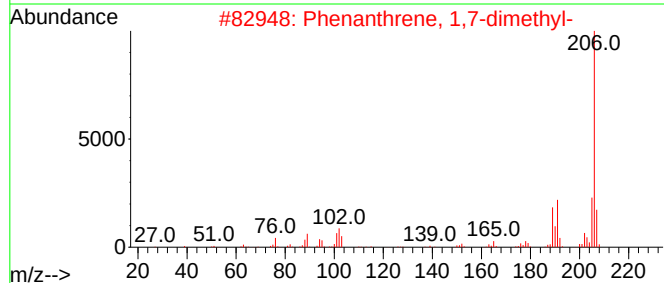
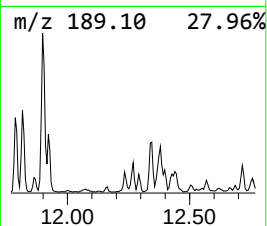
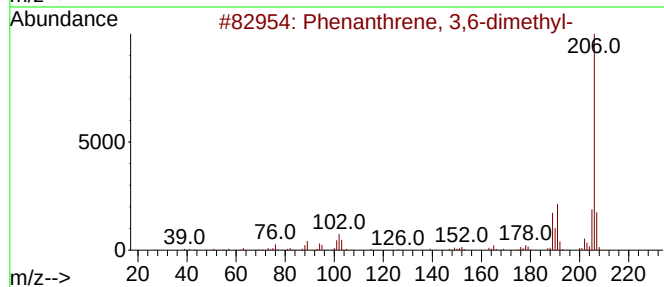
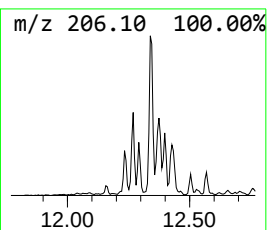
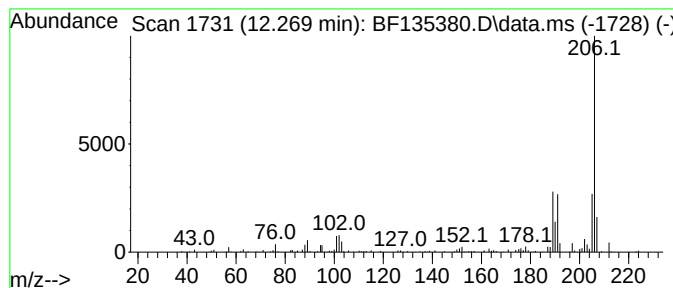
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Phenanthrene, 3,6-dimethyl- Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.269	2.31 ng	107688	Phenanthrene-d10	11.286

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 3,6-dimethyl-	206	C16H14	001576-67-6	94
2			Phenanthrene, 1,7-dimethyl-	206	C16H14	000483-87-4	94
3			di-p-Tolylacetylene	206	C16H14	002789-88-0	87
4			Phenanthrene, 2,7-dimethyl-	206	C16H14	001576-69-8	86
5			Phenanthrene, 2,3-dimethyl-	206	C16H14	003674-65-5	81



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF092123\
Data File : BF135380.D
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ClientSampleId :
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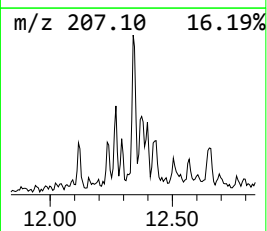
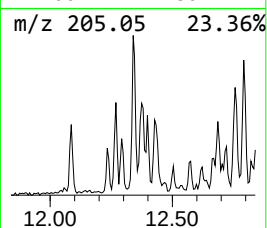
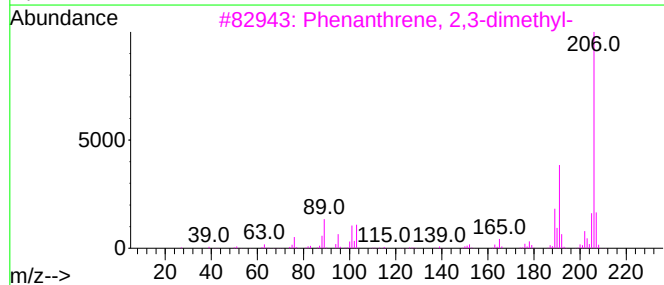
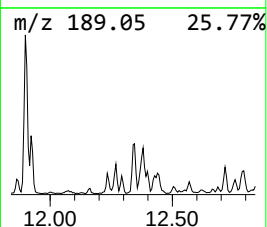
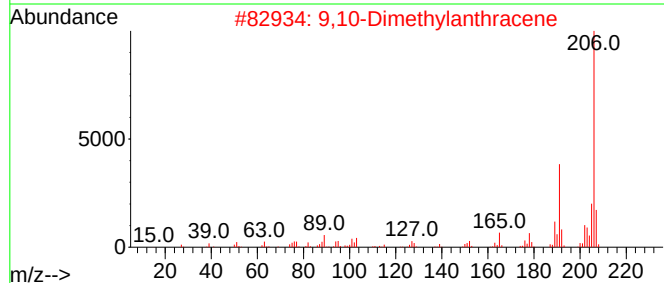
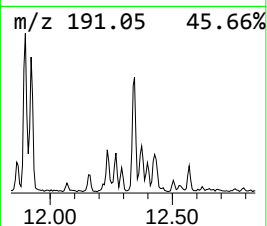
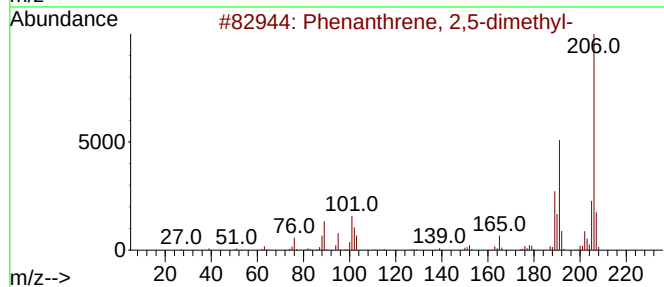
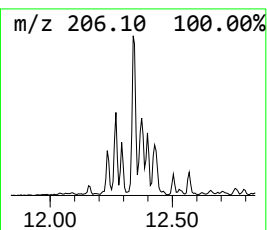
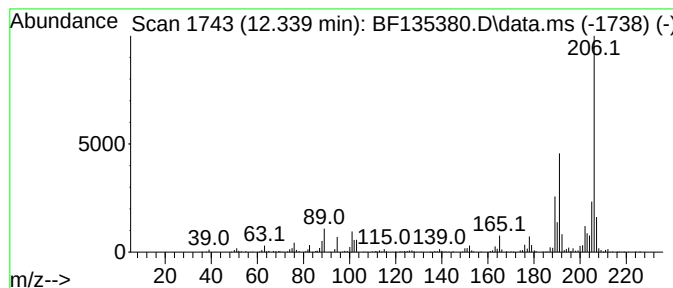
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Phenanthrene, 2,5-dimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.339	7.04 ng	328130	Phenanthrene-d10	11.286

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	97
2		9,10-Dimethylanthracene	206	C16H14	000781-43-1	96
3		Phenanthrene, 2,3-dimethyl-	206	C16H14	003674-65-5	94
4		Anthracene, 1,4-dimethyl-	206	C16H14	000781-92-0	94
5		Phenanthrene, 1,7-dimethyl-	206	C16H14	000483-87-4	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF092123\
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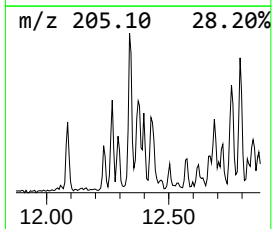
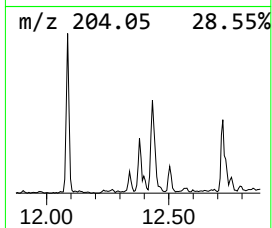
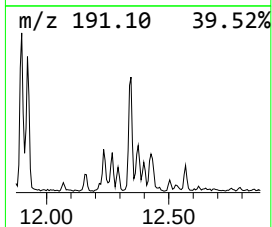
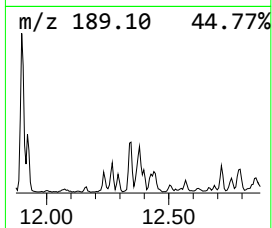
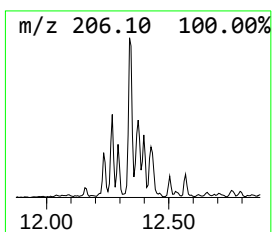
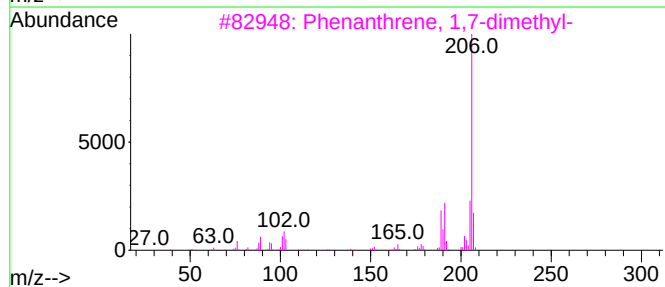
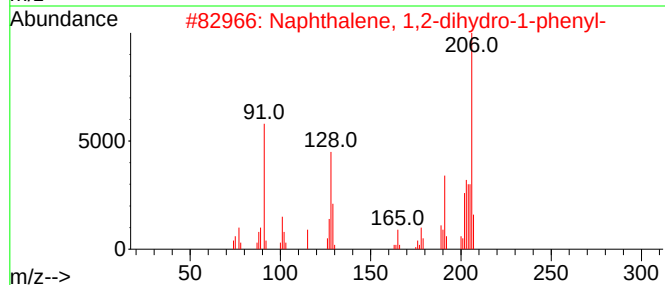
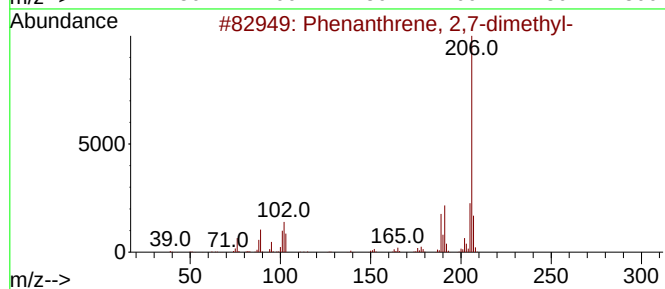
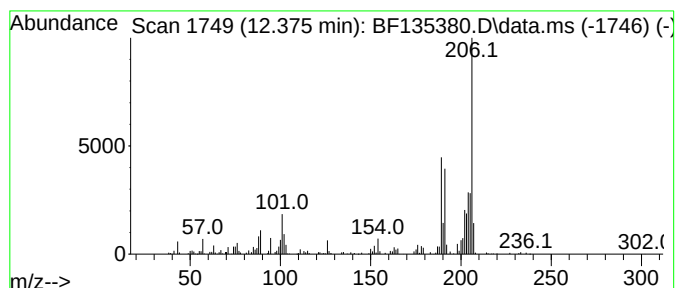
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Phenanthrene, 2,7-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.374	4.20 ng	195773	Phenanthrene-d10	11.286	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 2,7-dimethyl-	206	C16H14	001576-69-8	68
2	Naphthalene, 1,2-dihydro-1-phenyl-	206	C16H14	016606-46-5	60
3	Phenanthrene, 1,7-dimethyl-	206	C16H14	000483-87-4	55
4	9,10-Dimethylantracene	206	C16H14	000781-43-1	50
5	Anthracene, 1,4-dimethyl-	206	C16H14	000781-92-0	49



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF092123\
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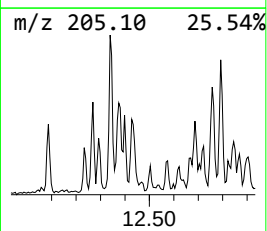
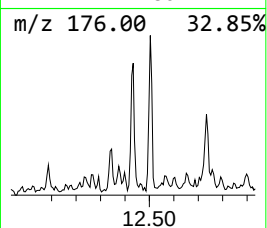
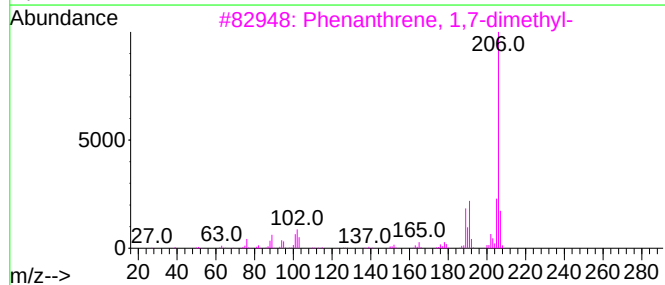
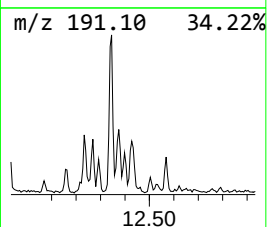
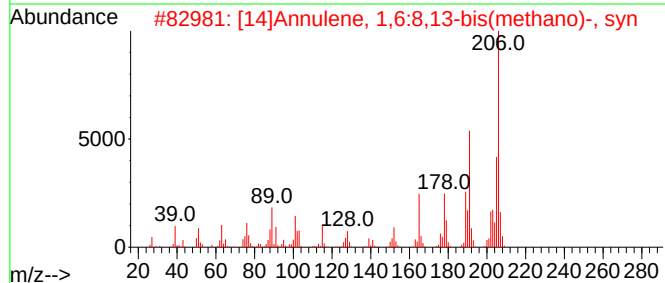
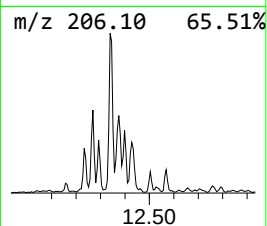
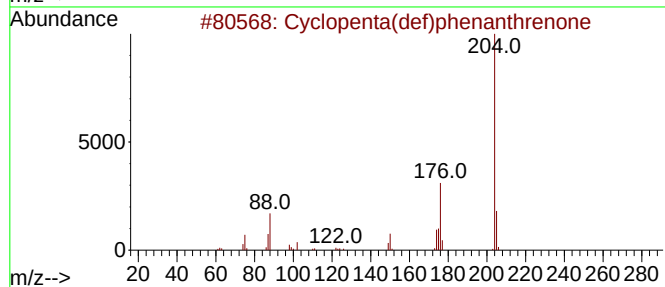
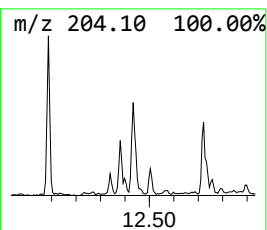
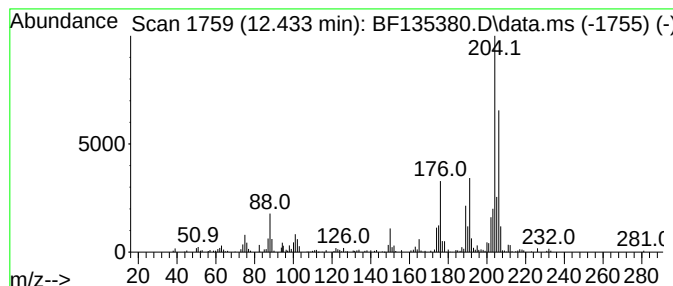
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Cyclopenta(def)phenanthrenone Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD		R.T.	
12.433	4.13 ng	192581	Phenanthrene-d10		11.286	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Cyclopenta(def)phenanthrenone		204	C15H8O	005737-13-3	94
2	[14]Annulene, 1,6:8,13-bis(metha...		206	C16H14	085385-68-8	46
3	Phenanthrene, 1,7-dimethyl-		206	C16H14	000483-87-4	46
4	1,2,4,8-Tetramethylbicyclo[6.3.0...		204	C15H24	137235-51-9	46
5	Phenanthrene, 3,6-dimethyl-		206	C16H14	001576-67-6	45



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF092123\
Data File : BF135380.D
Acq On : 21 Sep 2023 15:04
Operator : CG\JU
Sample : 04501-04
Misc :
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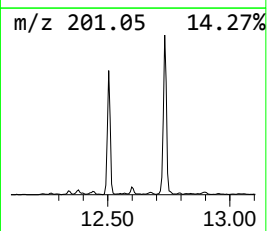
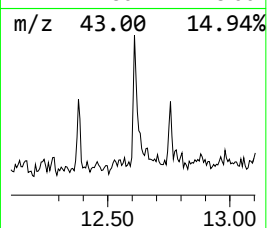
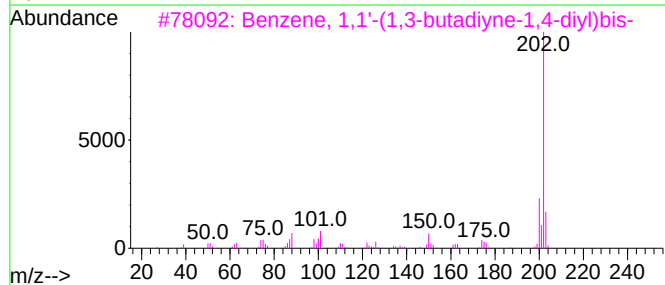
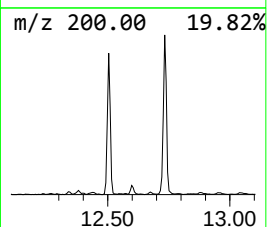
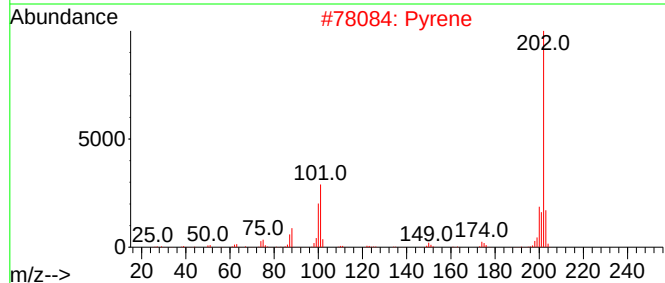
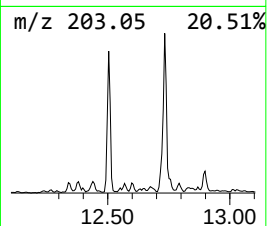
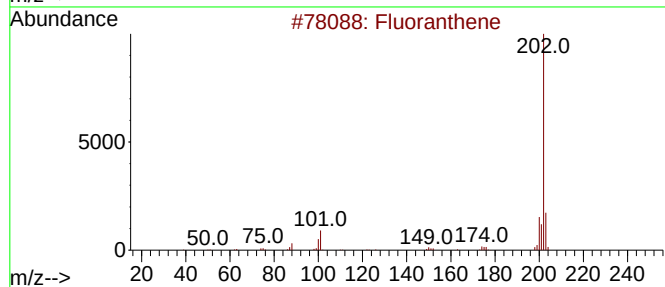
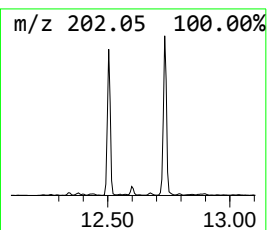
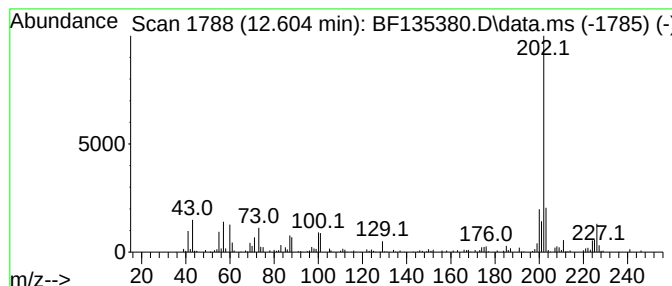
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Benzene, 1,1'-(1,3-butadiyn... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.604	2.66 ng	123825	Phenanthrene-d10	11.286	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Fluoranthene	202	C16H10	000206-44-0	70
2	Pyrene	202	C16H10	000129-00-0	58
3	Benzene, 1,1'-(1,3-butadiyne-1,4...	202	C16H10	000886-66-8	58
4	4,4'-Bis(tetrahydrothiopyran)	202	C10H18S2	116196-83-9	38
5	1,4-Pentadiyn-3-one, 1,5-diphenyl-	230	C17H10O	015814-30-9	37



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF092123\
Data File : BF135380.D
Acq On : 21 Sep 2023 15:04
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Sample : 04501-04
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Instrument :
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ClientSampleId :
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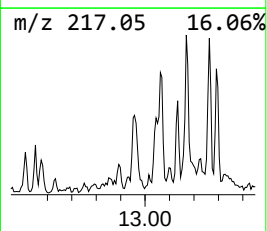
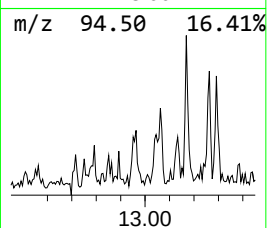
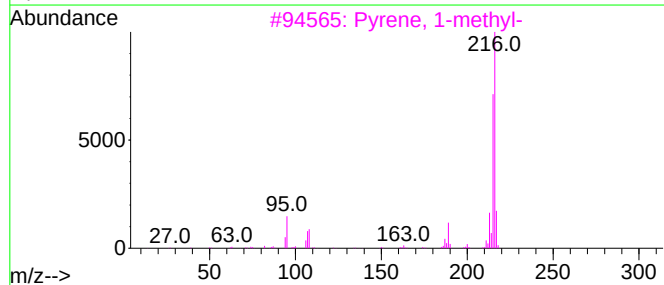
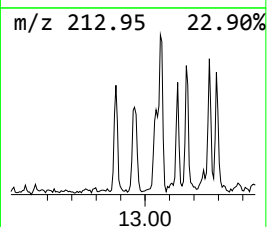
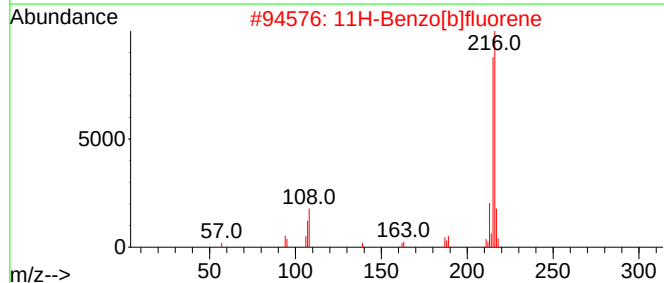
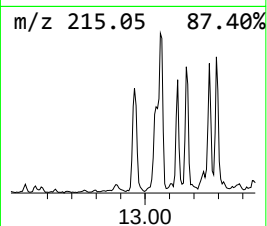
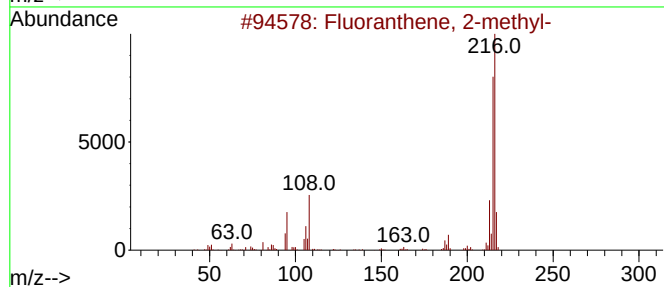
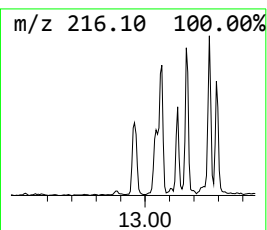
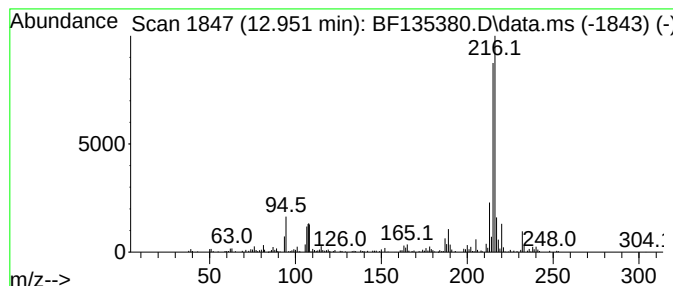
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Fluoranthene, 2-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.951	4.07 ng	273658	Chrysene-d12	13.945

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	95
2			11H-Benzo[b]fluorene	216	C17H12	000243-17-4	93
3			Pyrene, 1-methyl-	216	C17H12	002381-21-7	90
4			Pyrene, 2-methyl-	216	C17H12	003442-78-2	86
5			7H-Benzo[c]fluorene	216	C17H12	000205-12-9	81



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF092123\
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Sample : 04501-04
Misc :
ALS Vial : 12 Sample Multiplier: 1

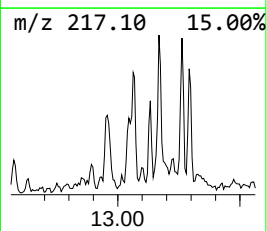
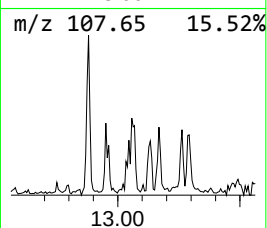
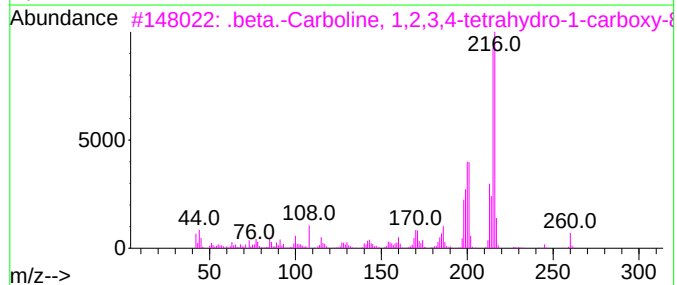
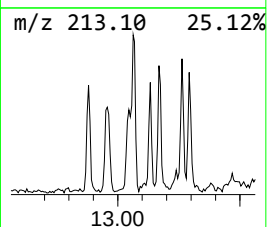
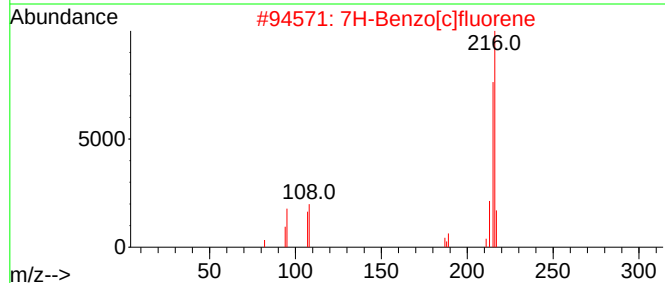
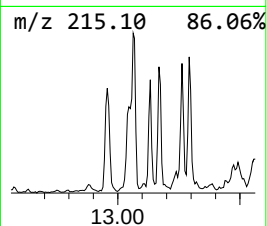
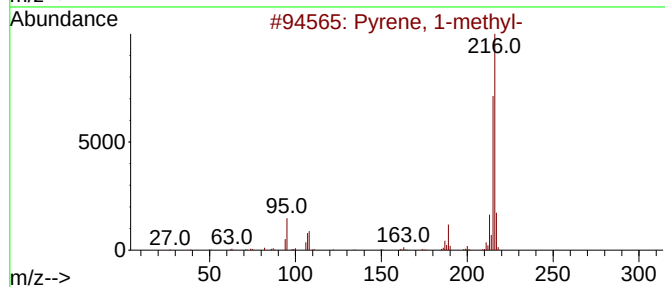
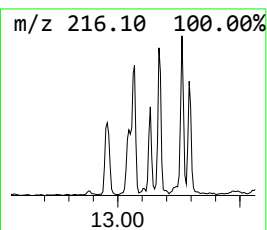
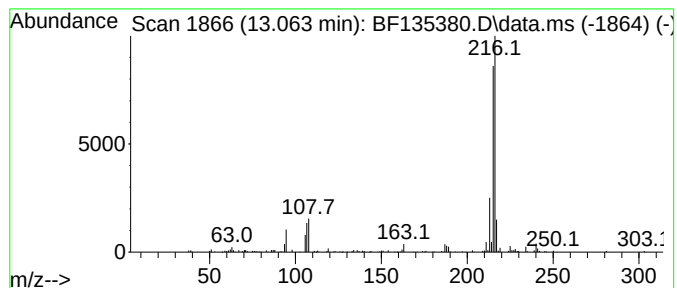
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF091123.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Pyrene, 1-methyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.063	3.29 ng	221396	Chrysene-d12	13.945	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pyrene, 1-methyl-	216	C17H12	002381-21-7	80
2	7H-Benzo[c]fluorene	216	C17H12	000205-12-9	74
3	.beta.-Carboline, 1,2,3,4-tetrahydro-1-carboxylic acid	260	C14H16N2O3	1000124-14-2	72
4	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	72
5	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF092123\
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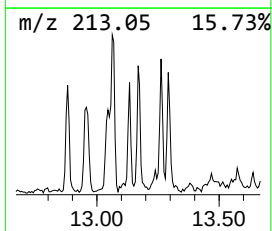
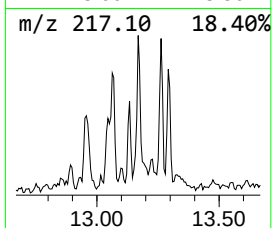
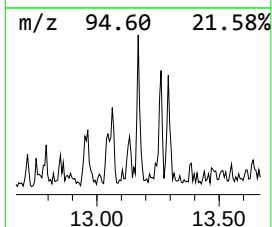
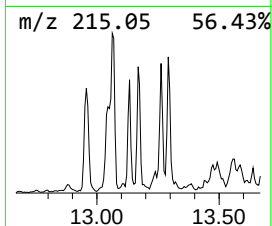
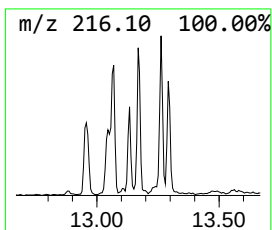
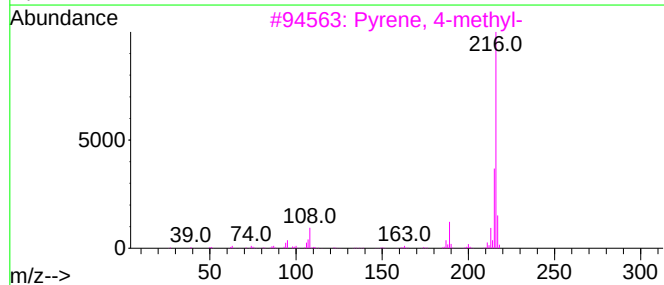
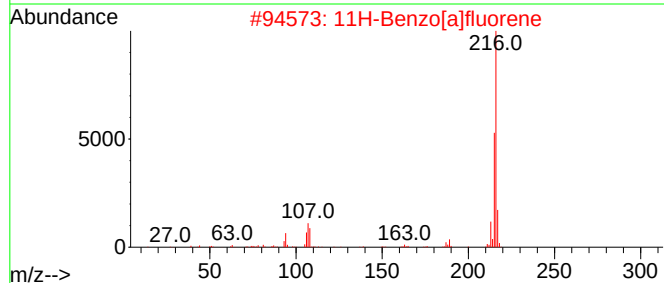
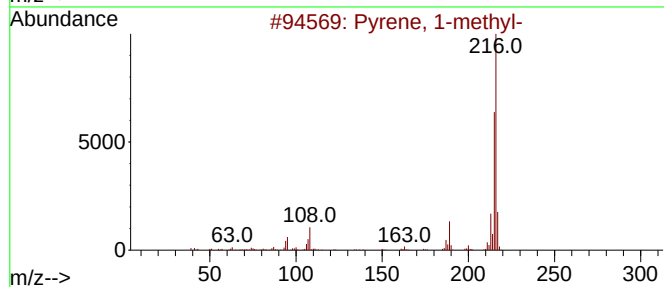
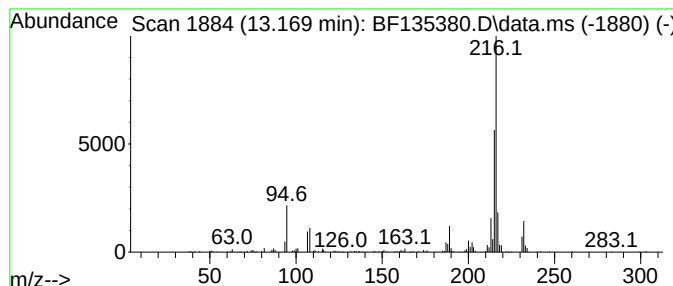
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 11H-Benzo[a]fluorene Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.169	3.27 ng	220337	Chrysene-d12	13.945	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pyrene, 1-methyl-	216	C17H12	002381-21-7	95
2	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	93
3	Pyrene, 4-methyl-	216	C17H12	003353-12-6	92
4	Pyrene, 2-methyl-	216	C17H12	003442-78-2	91
5	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF092123\
Data File : BF135380.D
Acq On : 21 Sep 2023 15:04
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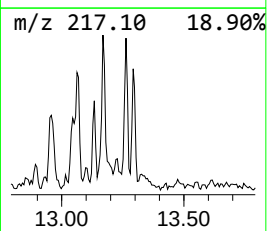
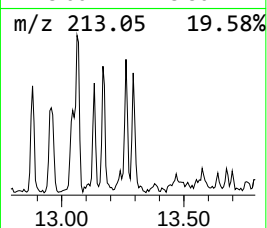
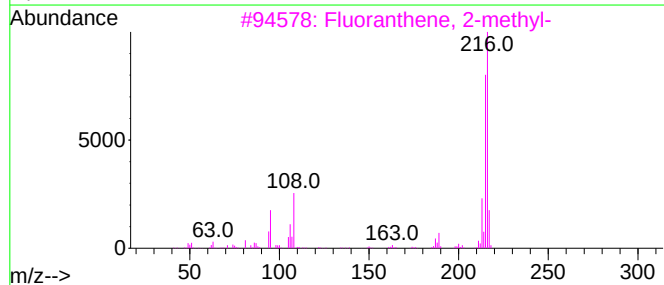
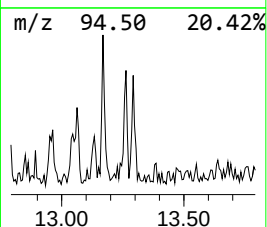
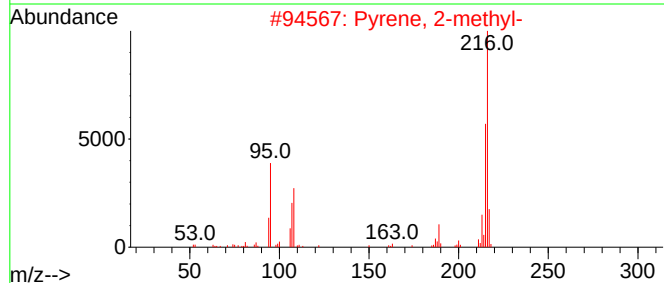
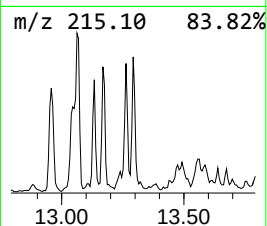
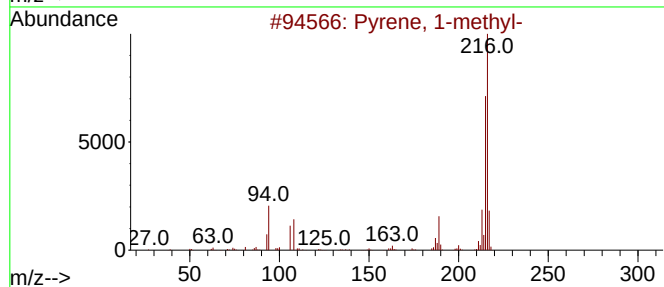
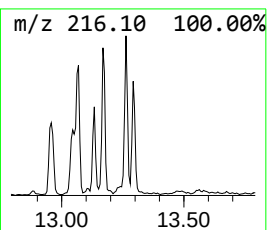
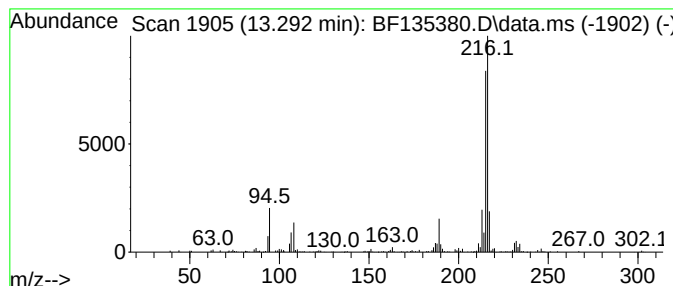
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Pyrene, 2-methyl- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.292	2.42 ng	162772	Chrysene-d12	13.945

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pyrene, 1-methyl-	216	C17H12	002381-21-7	97
2			Pyrene, 2-methyl-	216	C17H12	003442-78-2	93
3			Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	90
4			11H-Benzo[a]fluorene	216	C17H12	000238-84-6	87
5			11H-Benzo[b]fluorene	216	C17H12	000243-17-4	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF092123\
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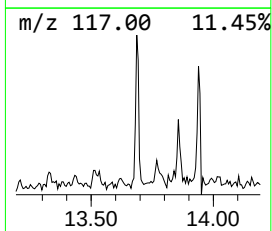
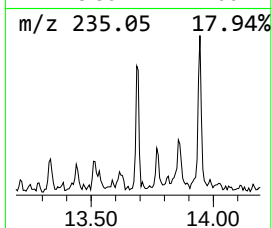
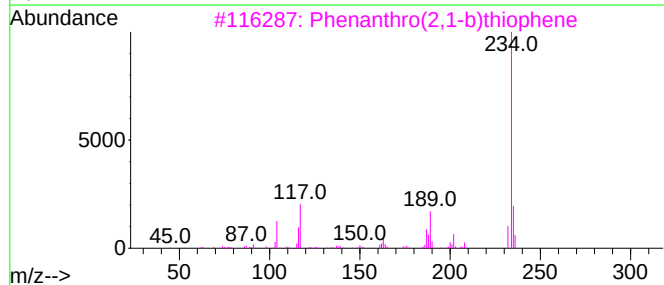
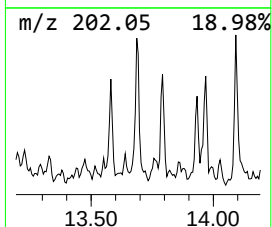
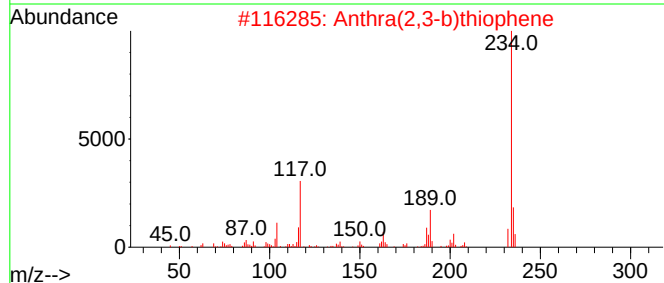
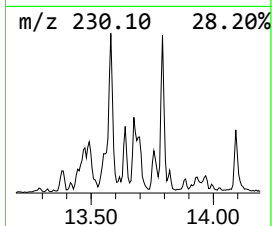
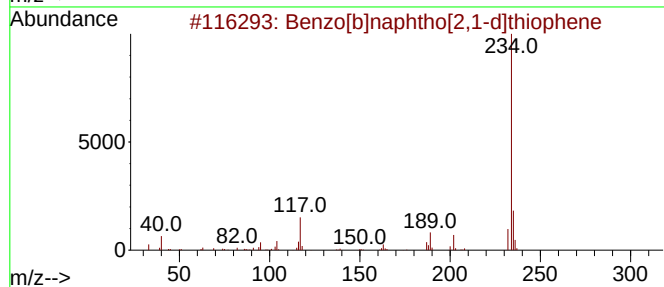
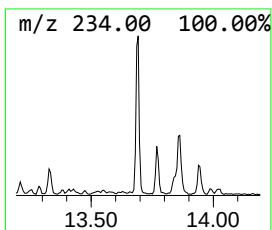
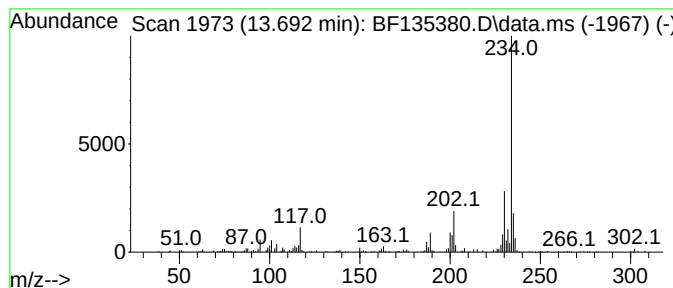
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 Benzo[b]naphtho[2,1-d]thiop... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.692	3.80 ng	255594	Chrysene-d12	13.945	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzo[b]naphtho[2,1-d]thiophene	234	C16H10S	000239-35-0	95
2	Anthra(2,3-b)thiophene	234	C16H10S	022108-55-0	76
3	Phenanthro(2,1-b)thiophene	234	C16H10S	000219-25-0	70
4	Benzo[b]naphtho[2,3-d]thiophene	234	C16H10S	000243-46-9	64
5	Benzo[b]naphtho[1,2-d]thiophene	234	C16H10S	000205-43-6	62



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF092123\
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Operator : CG\JU
Sample : 04501-04
Misc :
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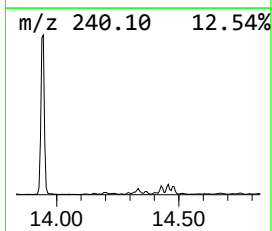
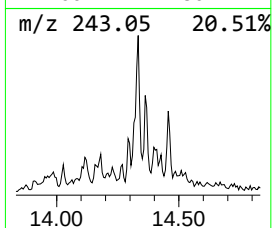
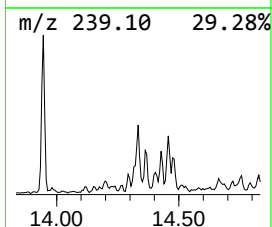
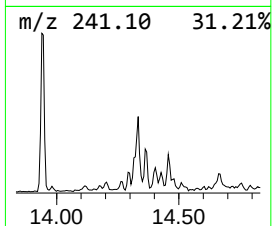
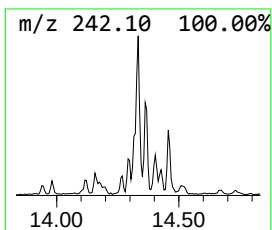
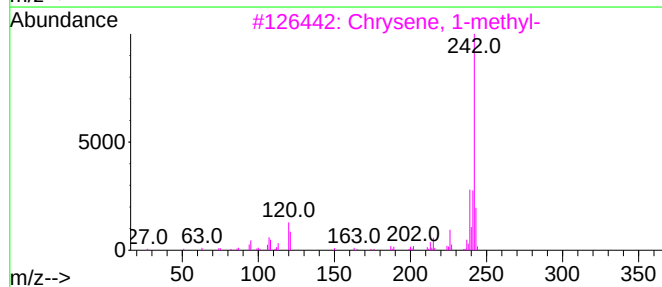
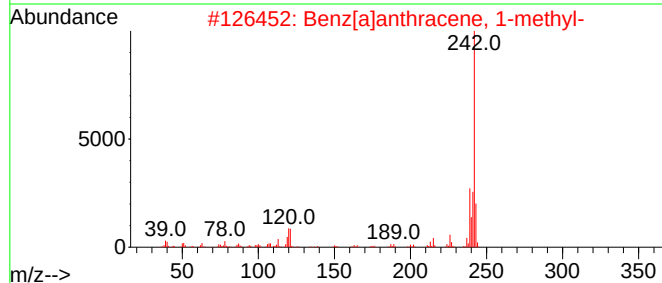
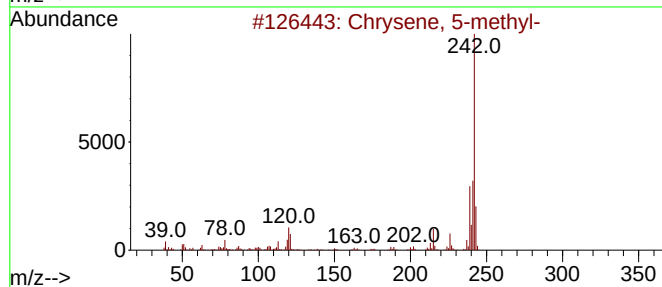
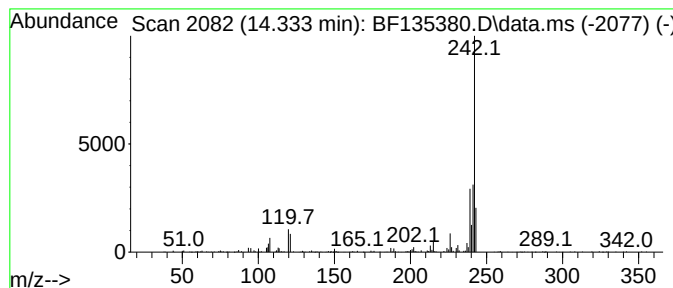
Instrument :
BNA_F
ClientSampleId :
TP-3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF091123.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 Chrysene, 5-methyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD		R.T.	
14.333	3.34 ng	224507	Chrysene-d12		13.945	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Chrysene, 5-methyl-		242	C19H14	003697-24-3	96
2	Benz[a]anthracene, 1-methyl-		242	C19H14	002498-77-3	95
3	Chrysene, 1-methyl-		242	C19H14	003351-28-8	93
4	Benz[a]anthracene, 8-methyl-		242	C19H14	002381-31-9	93
5	2-Methylchrysene		242	C19H14	003351-32-4	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF092123\
Data File : BF135380.D
Acq On : 21 Sep 2023 15:04
Operator : CG\JU
Sample : 04501-04
Misc :
ALS Vial : 12 Sample Multiplier: 1

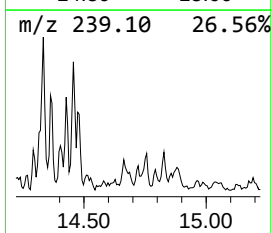
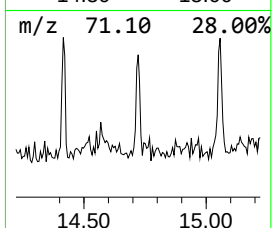
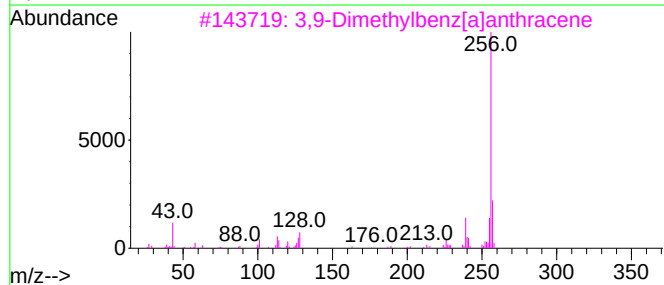
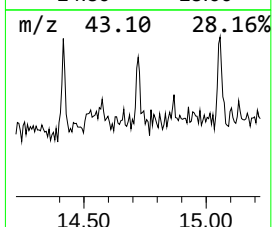
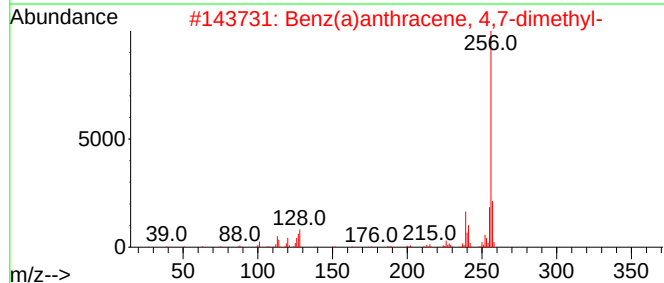
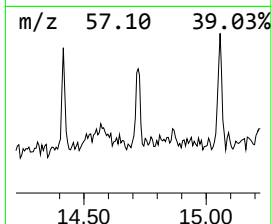
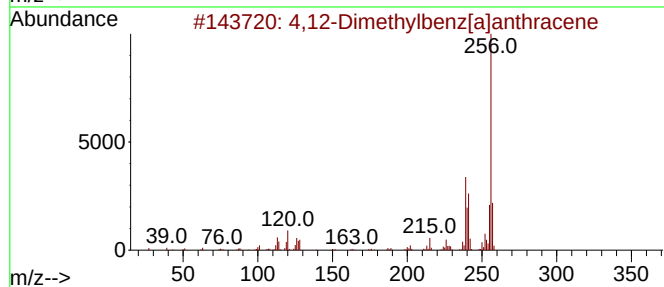
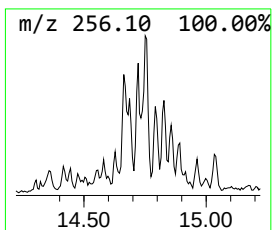
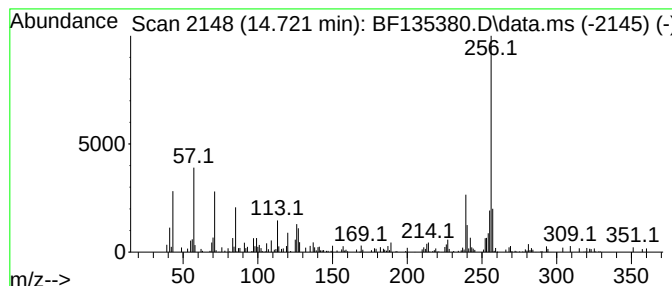
Instrument :
BNA_F
ClientSampleId :
TP-3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF091123.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 4,12-Dimethylbenz[a]anthracene Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.722	2.38 ng	73253	Perylene-d12	15.410	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4,12-Dimethylbenz[a]anthracene	256	C20H16	035187-19-0	90
2	Benz(a)anthracene, 4,7-dimethyl-	256	C20H16	035187-18-9	90
3	3,9-Dimethylbenz[a]anthracene	256	C20H16	000316-51-8	72
4	7,12-Dimethylbenz[a]anthracene	256	C20H16	000057-97-6	70
5	Benzo[c]phenanthrene, 5,8-dimethyl-	256	C20H16	054986-63-9	70



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF092123\
Data File : BF135380.D
Acq On : 21 Sep 2023 15:04
Operator : CG\JU
Sample : 04501-04
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TP-3

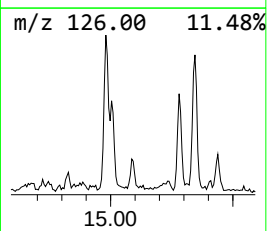
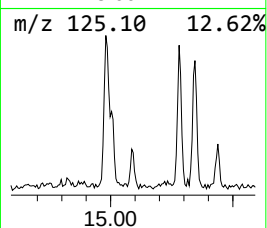
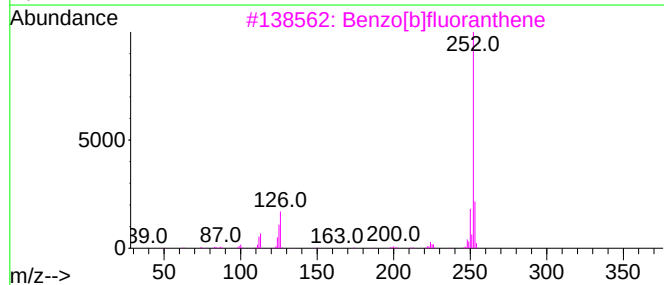
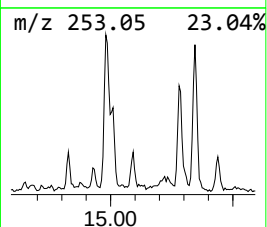
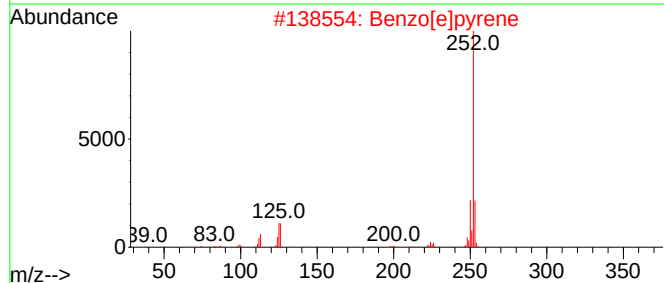
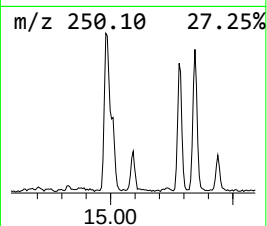
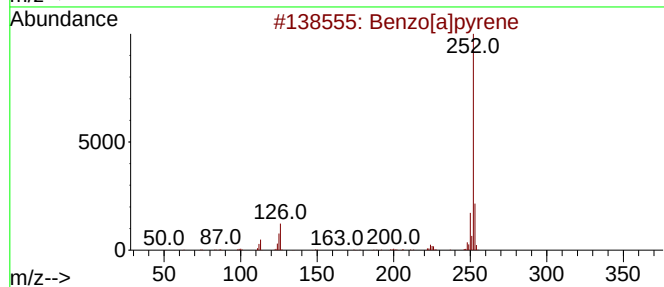
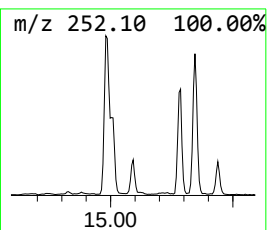
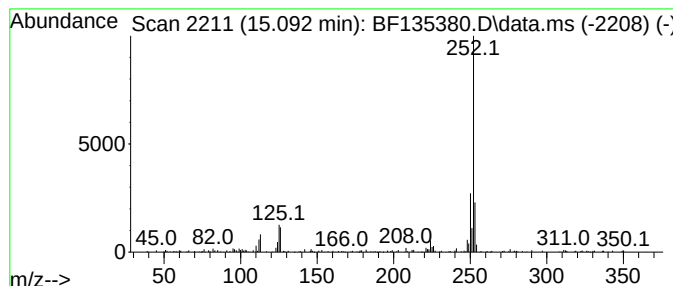
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 19 Benzo[e]pyrene Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.092	2.30 ng	71009	Perylene-d12	15.410

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzo[a]pyrene	252	C20H12	000050-32-8	97
2		Benzo[e]pyrene	252	C20H12	000192-97-2	94
3		Benzo[b]fluoranthene	252	C20H12	000205-99-2	93
4		Benzo[k]fluoranthene	252	C20H12	000207-08-9	91
5		Perylene	252	C20H12	000198-55-0	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF092123\
Data File : BF135380.D
Acq On : 21 Sep 2023 15:04
Operator : CG\JU
Sample : 04501-04
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TP-3

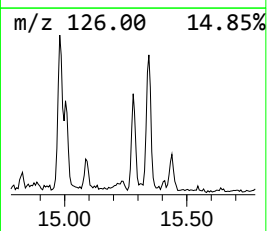
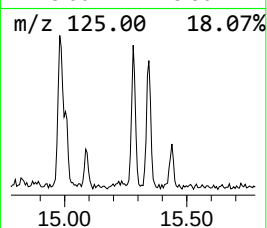
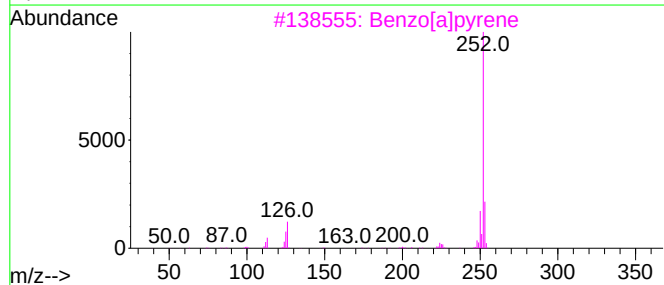
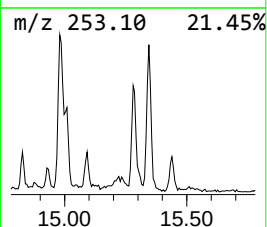
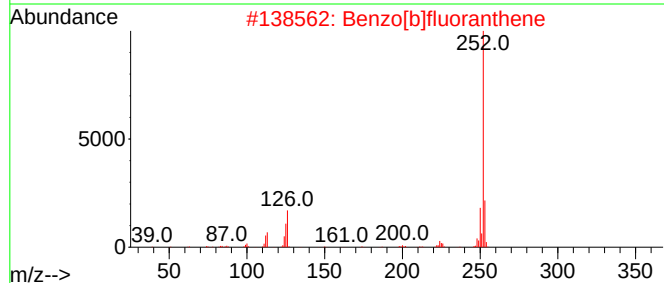
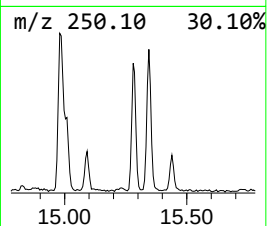
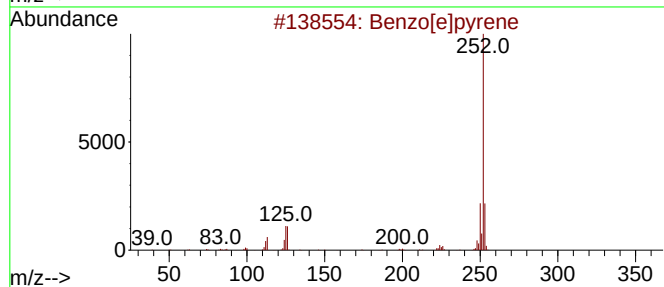
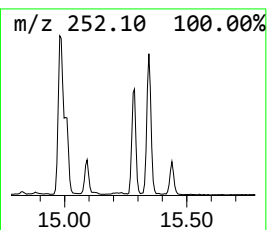
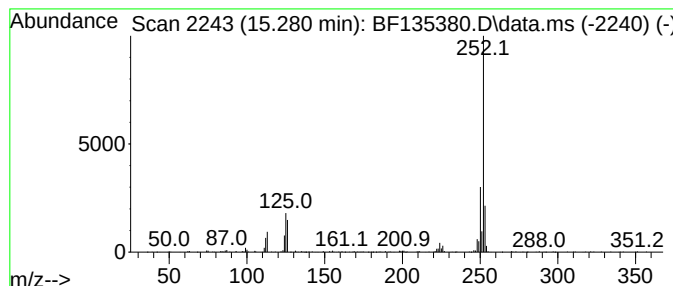
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF091123.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 20 Perylene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.280	9.20 ng	283652	Perylene-d12	15.410

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzo[e]pyrene	252	C20H12	000192-97-2	98
2		Benzo[b]fluoranthene	252	C20H12	000205-99-2	98
3		Benzo[a]pyrene	252	C20H12	000050-32-8	96
4		Perylene	252	C20H12	000198-55-0	93
5		Benzo[k]fluoranthene	252	C20H12	000207-08-9	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF092123\
 Data File : BF135380.D
 Acq On : 21 Sep 2023 15:04
 Operator : CG\JU
 Sample : 04501-04
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF091123.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
Phenanthrene, 2...	11.786	6.5	ng	303831	4	11.286	931629	20.0
Phenanthrene, 1...	11.816	11.7	ng	544796	4	11.286	931629	20.0
Anthracene, 2-m...	11.898	8.6	ng	398075	4	11.286	931629	20.0
1H-Cyclopropa[1...	11.922	4.7	ng	219991	4	11.286	931629	20.0
Naphthalene, 2-...	12.086	3.3	ng	155147	4	11.286	931629	20.0
9,10-Anthracene...	12.116	3.8	ng	174939	4	11.286	931629	20.0
Phenanthrene, 3...	12.269	2.3	ng	107688	4	11.286	931629	20.0
Phenanthrene, 2...	12.339	7.0	ng	328130	4	11.286	931629	20.0
Phenanthrene, 2...	12.374	4.2	ng	195773	4	11.286	931629	20.0
Cyclopenta(def)...	12.433	4.1	ng	192581	4	11.286	931629	20.0
Benzene, 1,1'-(...	12.604	2.7	ng	123825	4	11.286	931629	20.0
Fluoranthene, 2...	12.951	4.1	ng	273658	5	13.945	1346310	20.0
Pyrene, 1-methyl-	13.063	3.3	ng	221396	5	13.945	1346310	20.0
11H-Benzo[a]flu...	13.169	3.3	ng	220337	5	13.945	1346310	20.0
Pyrene, 2-methyl-	13.292	2.4	ng	162772	5	13.945	1346310	20.0
Benzo[b]naphtho...	13.692	3.8	ng	255594	5	13.945	1346310	20.0
Chrysene, 5-met...	14.333	3.3	ng	224507	5	13.945	1346310	20.0
4,12-Dimethylbe...	14.722	2.4	ng	73253	6	15.410	616491	20.0
Benzo[e]pyrene	15.092	2.3	ng	71009	6	15.410	616491	20.0
Perylene	15.280	9.2	ng	283652	6	15.410	616491	20.0