

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF062524\
 Data File : BF138344.D
 Acq On : 25 Jun 2024 20:08
 Operator : CG\JU
 Sample : P3021-10 10X
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 ARS20-0015

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-BF061224.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF138344.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.146	5	10	17	rVB	382146	461469	19.69%	1.898%
2	3.957	313	318	324	rVB	18606	23545	1.00%	0.097%
3	5.498	576	580	590	rVB	403819	354261	15.11%	1.457%
4	6.504	748	751	768	rVB	425822	343318	14.65%	1.412%
5	6.887	812	816	819	rBV	1409900	1143713	48.80%	4.703%
6	7.439	907	910	915	rBV	256734	203785	8.69%	0.838%
7	8.169	1030	1034	1036	rBV	1571035	1365723	58.27%	5.616%
8	8.186	1036	1037	1041	rVB	54483	32554	1.39%	0.134%
9	8.975	1168	1171	1176	rVB	29302	24230	1.03%	0.100%
10	9.239	1212	1216	1219	rBV	528087	409675	17.48%	1.685%
11	9.581	1270	1274	1276	rBV	45341	44664	1.91%	0.184%
12	9.781	1305	1308	1311	rVB	54244	43118	1.84%	0.177%
13	9.922	1328	1332	1335	rBV	1655331	1316138	56.15%	5.412%
14	9.951	1335	1337	1341	rVB	85913	70522	3.01%	0.290%
15	10.016	1344	1348	1353	rVB2	66174	86997	3.71%	0.358%
16	10.122	1362	1366	1370	rVV	52692	47819	2.04%	0.197%
17	10.163	1370	1373	1376	rVB	27714	25440	1.09%	0.105%
18	10.233	1382	1385	1388	rBV	29378	28145	1.20%	0.116%
19	10.328	1396	1401	1403	rBV2	30727	43040	1.84%	0.177%
20	10.463	1421	1424	1430	rVB2	72199	82625	3.53%	0.340%
21	10.533	1430	1436	1437	rBV2	24573	32303	1.38%	0.133%
22	10.628	1446	1452	1457	rVB3	43125	71976	3.07%	0.296%
23	10.704	1462	1465	1469	rVV	236375	199437	8.51%	0.820%
24	10.833	1482	1487	1491	rBV4	36154	55563	2.37%	0.228%
25	11.016	1512	1518	1520	rBV3	43430	55050	2.35%	0.226%
26	11.045	1520	1523	1525	rVB2	26254	24014	1.02%	0.099%
27	11.139	1535	1539	1541	rBV3	31747	40116	1.71%	0.165%
28	11.210	1548	1551	1555	rVB2	40700	37113	1.58%	0.153%
29	11.269	1555	1561	1564	rVV2	47581	74955	3.20%	0.308%
30	11.304	1564	1567	1574	rVB	67745	66342	2.83%	0.273%
31	11.404	1581	1584	1587	rBV	1343594	1331711	56.82%	5.476%
32	11.433	1587	1589	1592	rVV	962484	776413	33.13%	3.193%
33	11.480	1595	1597	1600	rVV	248995	187544	8.00%	0.771%
34	11.510	1600	1602	1606	rVV2	28431	30806	1.31%	0.127%
35	11.633	1617	1623	1626	rBV2	51266	63592	2.71%	0.261%
36	11.704	1631	1635	1638	rVV2	52187	53896	2.30%	0.222%

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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-BF061224.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

37	11.739	1638	1641	1646	rVB2	48320	50142	2.14%	0.206%
38	11.822	1652	1655	1659	rVB	26161	29028	1.24%	0.119%
39	11.910	1664	1670	1672	rBV	211371	195945	8.36%	0.806%
40	11.933	1672	1674	1677	rVV	268944	234683	10.01%	0.965%
41	11.980	1677	1682	1685	rVV2	93458	101370	4.32%	0.417%
42	12.016	1685	1688	1691	rVV2	287270	338343	14.44%	1.391%
43	12.039	1691	1692	1697	rVB	162878	125285	5.35%	0.515%
44	12.110	1697	1704	1707	rBV6	25065	56015	2.39%	0.230%
45	12.204	1717	1720	1722	rBV	158657	132186	5.64%	0.544%
46	12.227	1722	1724	1727	rVB2	90107	76031	3.24%	0.313%
47	12.351	1740	1745	1748	rBV3	71087	83933	3.58%	0.345%
48	12.380	1748	1750	1752	rBV	48989	35352	1.51%	0.145%
49	12.404	1752	1754	1757	rVB	47658	39646	1.69%	0.163%
50	12.457	1759	1763	1766	rBV	156570	165406	7.06%	0.680%
51	12.492	1766	1769	1771	rVV2	91717	100091	4.27%	0.412%
52	12.516	1771	1773	1775	rVB	52997	37124	1.58%	0.153%
53	12.539	1775	1777	1783	rBV2	115029	135686	5.79%	0.558%
54	12.622	1787	1791	1794	rBV	1935486	1625049	69.33%	6.682%
55	12.663	1794	1798	1800	rBV	33644	32169	1.37%	0.132%
56	12.680	1800	1801	1804	rVB2	42147	31813	1.36%	0.131%
57	12.716	1804	1807	1809	rVB	50540	44618	1.90%	0.183%
58	12.745	1809	1812	1814	rBV2	45067	51489	2.20%	0.212%
59	12.769	1814	1816	1819	rVB	67075	60094	2.56%	0.247%
60	12.851	1823	1830	1833	rBV	1928245	1898387	80.99%	7.806%
61	12.898	1835	1838	1843	rVB2	79768	100444	4.29%	0.413%
62	12.969	1846	1850	1851	rBV	93883	92452	3.94%	0.380%
63	12.986	1851	1853	1860	rVB2	507947	509021	21.72%	2.093%
64	13.063	1861	1866	1870	rBV3	149595	238359	10.17%	0.980%
65	13.121	1874	1876	1878	rBV3	29363	30601	1.31%	0.126%
66	13.151	1878	1881	1882	rBV2	113046	104842	4.47%	0.431%
67	13.174	1882	1885	1888	rVB	261672	259693	11.08%	1.068%
68	13.221	1888	1893	1894	rBV3	38498	53938	2.30%	0.222%
69	13.239	1894	1896	1899	rBV	144651	130385	5.56%	0.536%
70	13.280	1899	1903	1906	rBV2	215432	222204	9.48%	0.914%
71	13.304	1906	1907	1910	rVB	42366	34659	1.48%	0.143%
72	13.339	1910	1913	1916	rBV3	38190	39388	1.68%	0.162%
73	13.374	1916	1919	1921	rBV	118136	106733	4.55%	0.439%
74	13.404	1921	1924	1927	rBV2	139725	120642	5.15%	0.496%
75	13.563	1947	1951	1952	rBV2	47832	63461	2.71%	0.261%
76	13.680	1968	1971	1976	rVB	141952	144578	6.17%	0.595%

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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-BF061224.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

77	13.792	1986	1990	1993	rBV2	138365	187245	7.99%	0.770%
78	13.821	1993	1995	1997	rVV	169560	140751	6.01%	0.579%
79	13.845	1997	1999	2003	rVV2	135213	177817	7.59%	0.731%
80	13.886	2004	2006	2014	rVB	140540	167265	7.14%	0.688%
81	14.045	2025	2033	2036	rBV2	1696524	2343872	100.00%	9.638%
82	14.068	2036	2037	2044	rVB	772219	676291	28.85%	2.781%
83	14.151	2049	2051	2053	rVV	94436	71589	3.05%	0.294%
84	14.392	2090	2092	2094	rBV2	53172	56782	2.42%	0.233%
85	14.433	2096	2099	2102	rVV	162546	177856	7.59%	0.731%
86	14.463	2102	2104	2108	rVV	103078	103919	4.43%	0.427%
87	14.557	2117	2120	2122	rBV	70615	66302	2.83%	0.273%
88	15.086	2206	2210	2214	rBV	508388	812169	34.65%	3.340%
89	15.198	2226	2229	2234	rVB	110134	124338	5.30%	0.511%
90	15.392	2259	2262	2268	rVB	294342	366020	15.62%	1.505%
91	15.457	2269	2273	2279	rBV	438914	522691	22.30%	2.149%
92	15.521	2280	2284	2287	rBV	682049	826494	35.26%	3.399%
93	16.998	2531	2535	2544	rVB2	162675	314308	13.41%	1.292%

Sum of corrected areas: 24318606

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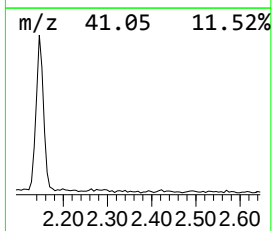
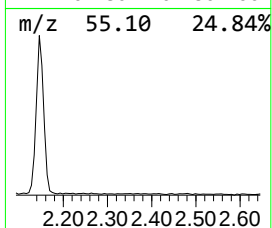
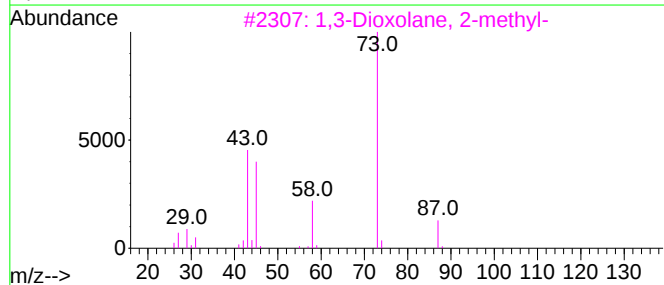
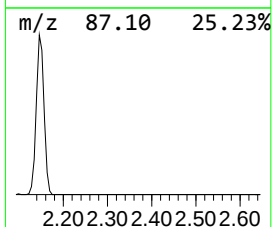
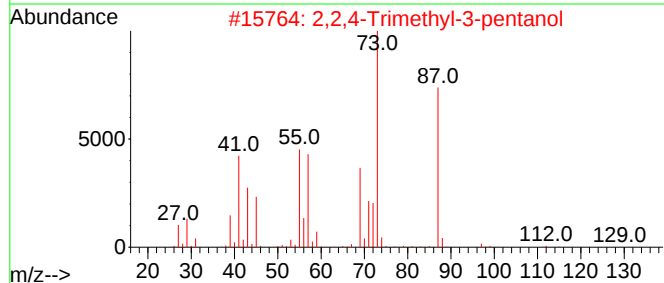
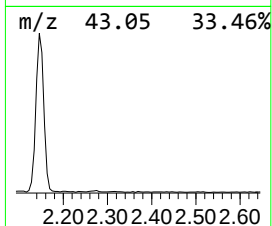
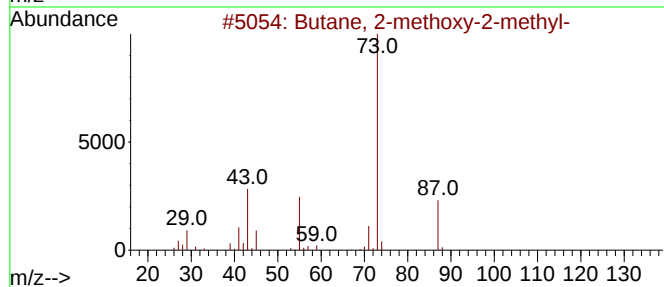
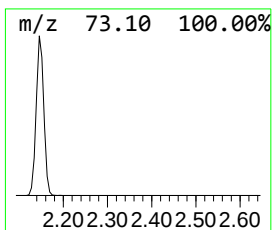
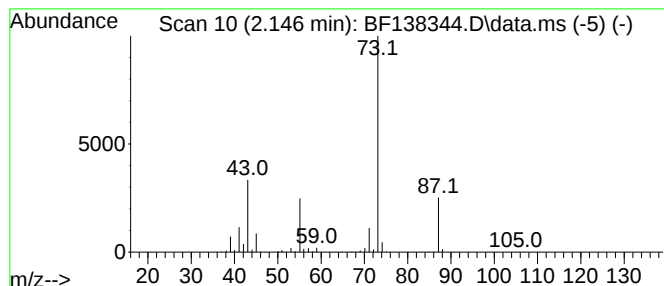
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF061224.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.146	8.07 ng	461469	1,4-Dichlorobenzene-d4	6.887

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
3			1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	25
4			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	23
5			Silane, tetramethyl-	88	C4H12Si	000075-76-3	17



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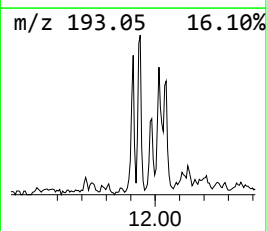
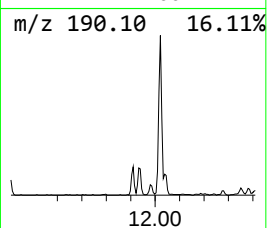
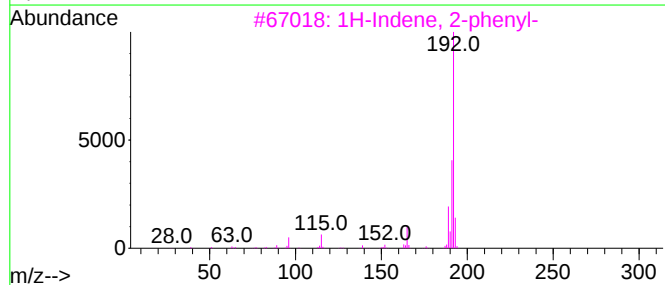
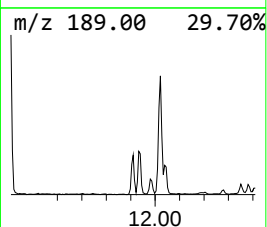
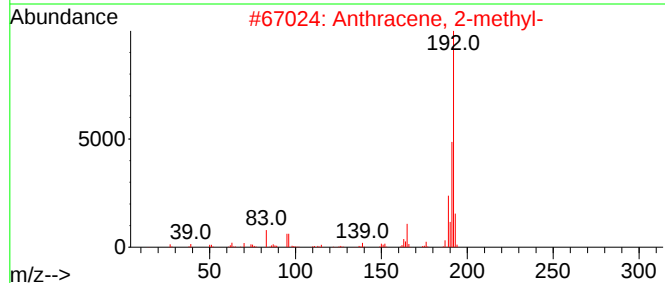
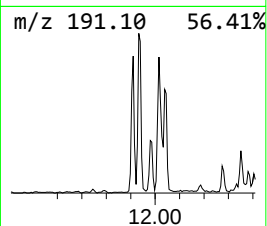
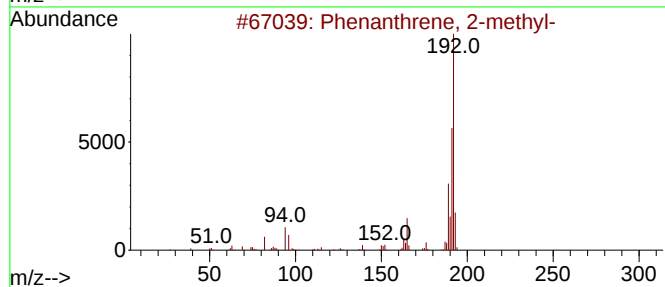
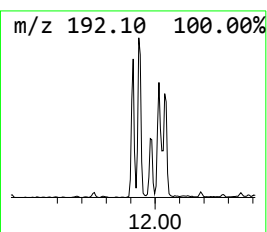
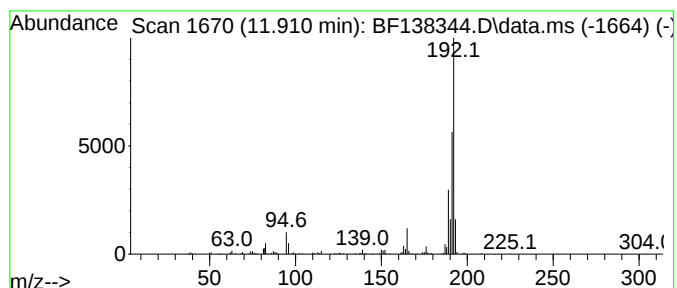
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF061224.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 Phenanthrene, 2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.910	2.94 ng	195945	Phenanthrene-d10	11.404

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	98
2		Anthracene, 2-methyl-	192	C15H12	000613-12-7	95
3		1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	93
4		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	93
5		Anthracene, 9-methyl-	192	C15H12	000779-02-2	91



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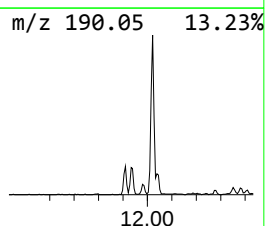
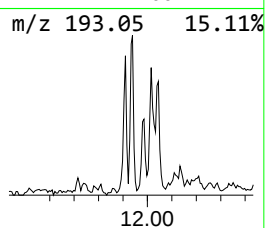
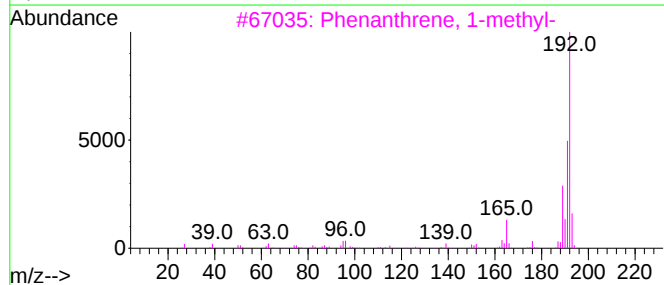
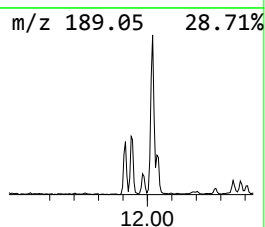
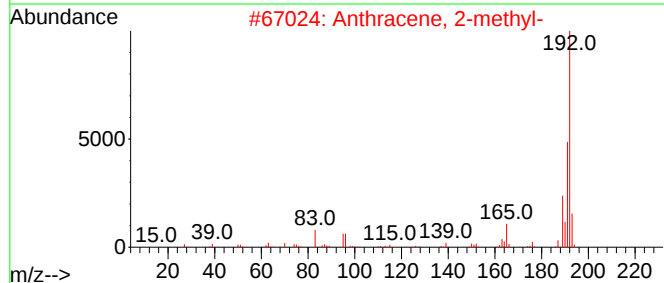
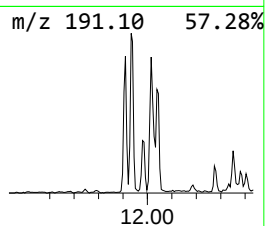
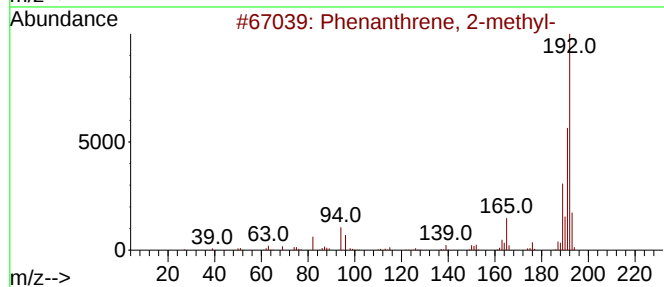
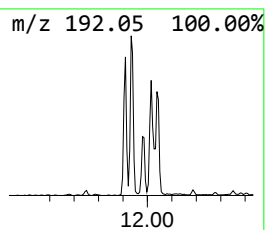
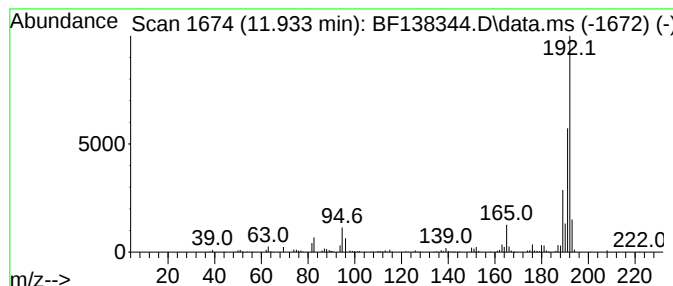
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF061224.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.933	3.52 ng	234683	Phenanthrene-d10	11.404

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	98
2			Anthracene, 2-methyl-	192	C15H12	000613-12-7	96
3			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	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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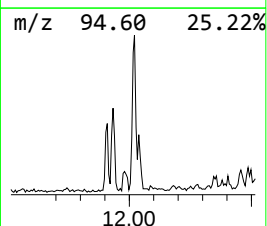
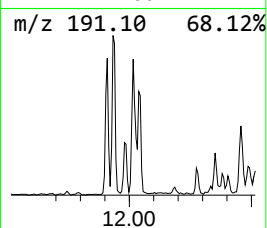
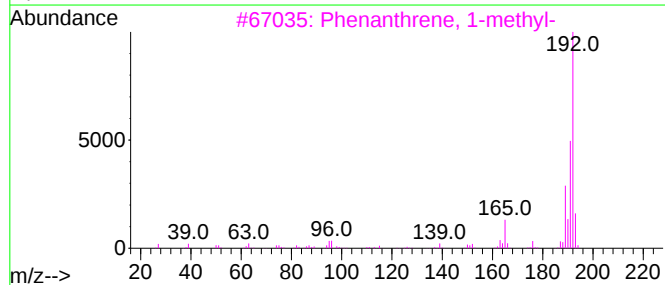
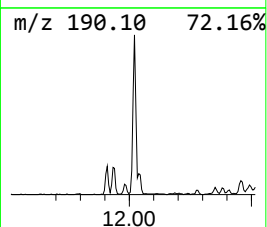
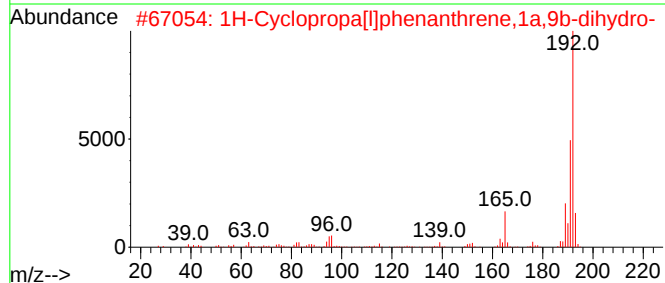
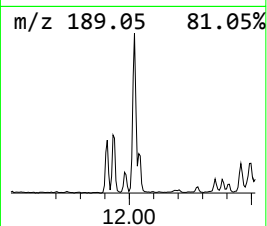
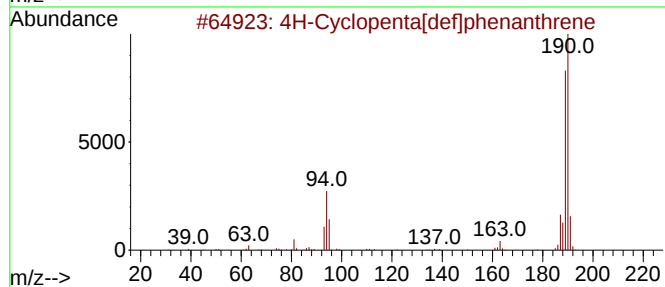
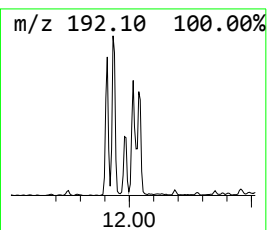
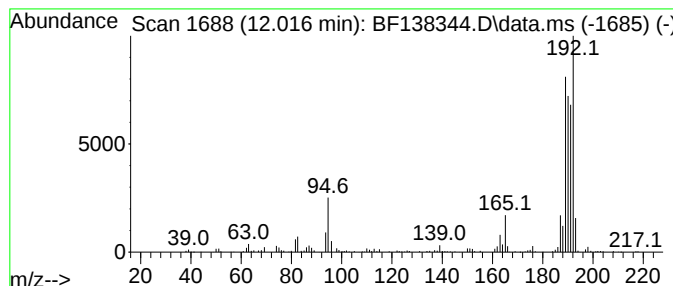
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.016	5.08 ng	338343	Phenanthrene-d10	11.404

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	55
2		1H-Cyclopropa[l]phenanthrene,1a,...	192	C15H12	000949-41-7	50
3		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	50
4		5H-Dibenzo[a,d]cycloheptene	192	C15H12	000256-81-5	50
5		Anthracene, 1-methyl-	192	C15H12	000610-48-0	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF062524\
Data File : BF138344.D
Acq On : 25 Jun 2024 20:08
Operator : CG\JU
Sample : P3021-10 10X
Misc :
ALS Vial : 21 Sample Multiplier: 1

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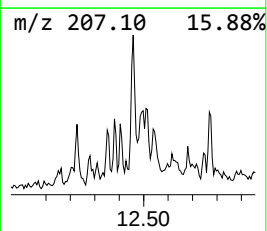
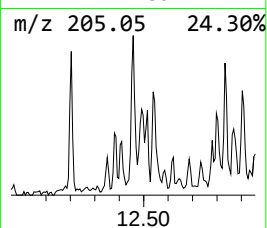
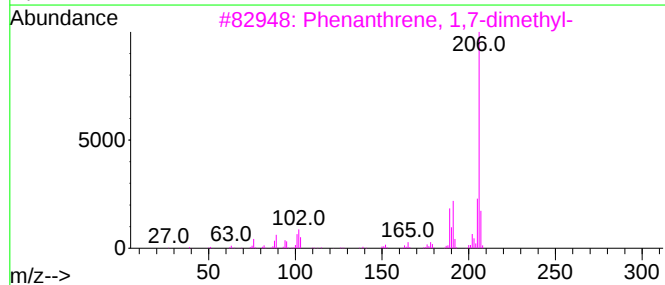
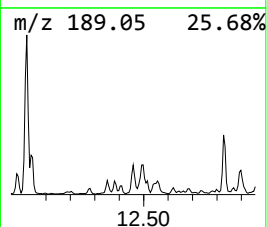
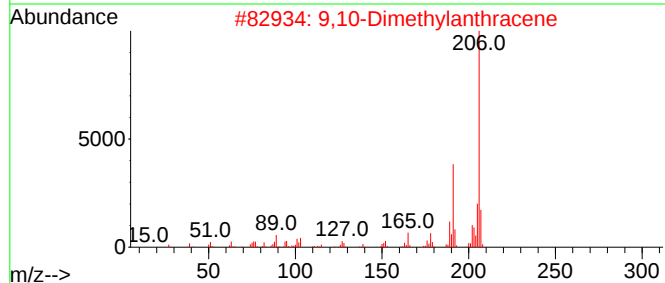
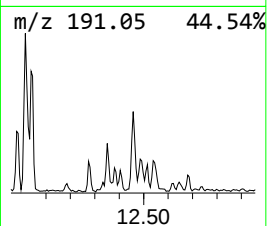
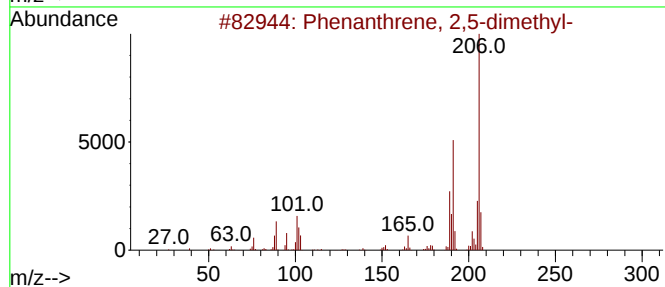
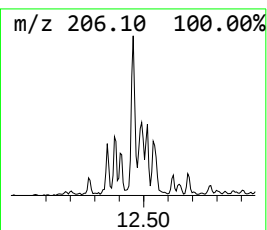
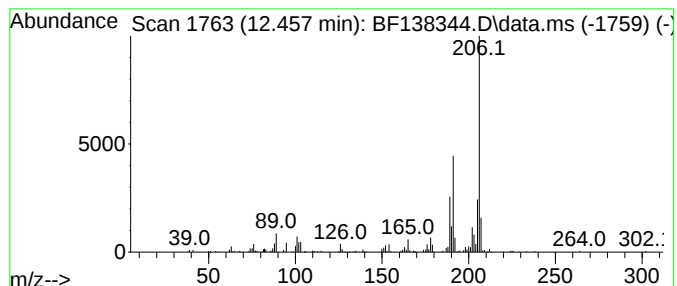
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Phenanthrene, 2,5-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.457	2.48 ng	165406	Phenanthrene-d10	11.404

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	93
2			9,10-Dimethylanthracene	206	C16H14	000781-43-1	90
3			Phenanthrene, 1,7-dimethyl-	206	C16H14	000483-87-4	89
4			Anthracene, 1,4-dimethyl-	206	C16H14	000781-92-0	89
5			Phenanthrene, 2,3-dimethyl-	206	C16H14	003674-65-5	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF062524\
Data File : BF138344.D
Acq On : 25 Jun 2024 20:08
Operator : CG\JU
Sample : P3021-10 10X
Misc :
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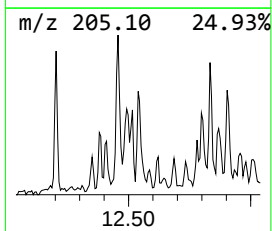
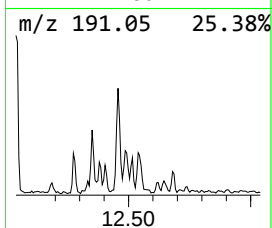
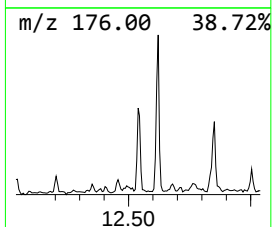
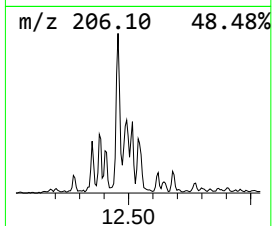
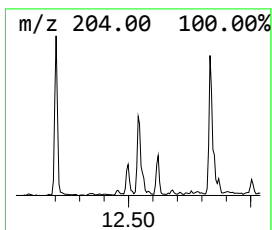
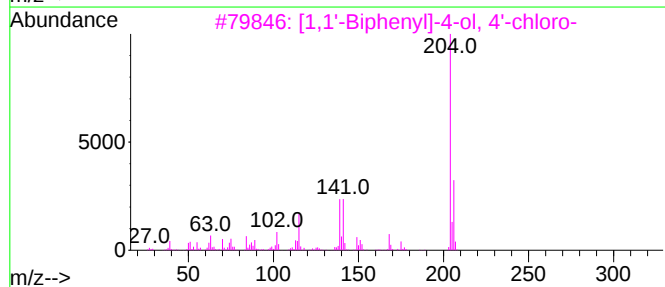
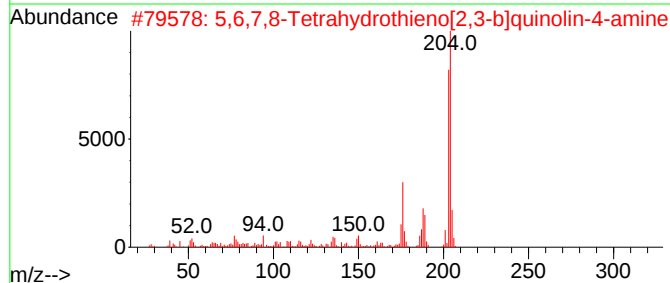
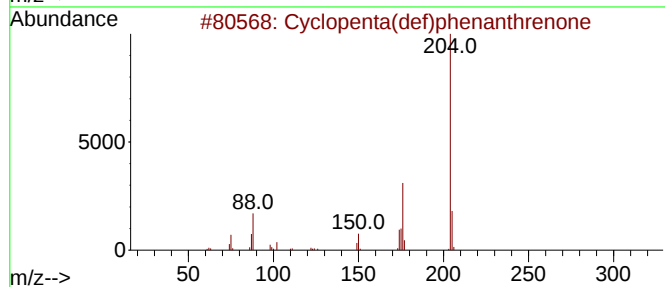
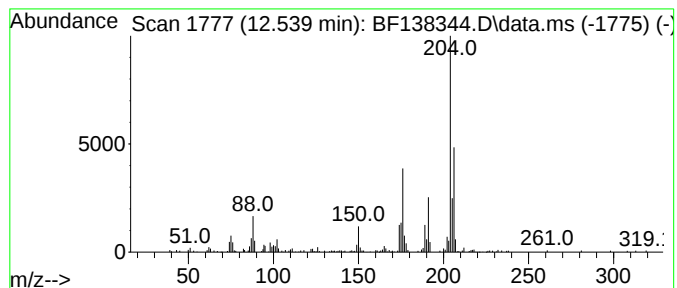
Instrument :
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Cyclopenta(def)phenanthrenone Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD		R.T.	
12.539	2.04 ng	135686	Phenanthrene-d10		11.404	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Cyclopenta(def)phenanthrenone		204	C15H8O	005737-13-3	93
2	5,6,7,8-Tetrahydrothieno[2,3-b]q...		204	C11H12N2S	122914-50-5	50
3	[1,1'-Biphenyl]-4-ol, 4'-chloro-		204	C12H9ClO	028034-99-3	43
4	2-Chloro-4-phenylphenol		204	C12H9ClO	000092-04-6	38
5	2,4-Dichloro-5-ethyl-3-methylphe...		290	C13H16Cl2O3	1000487-27-7	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF062524\
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Sample : P3021-10 10X
Misc :
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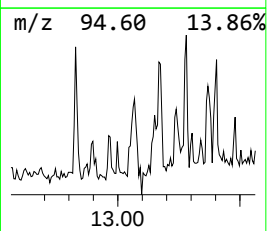
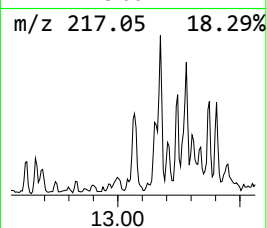
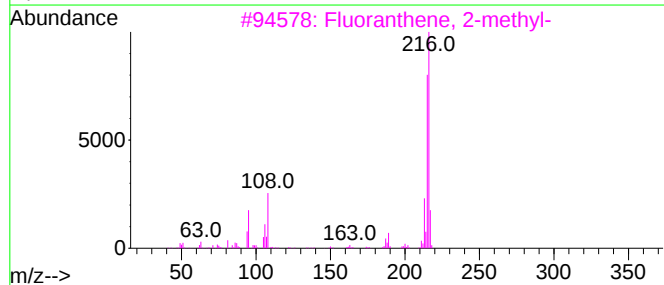
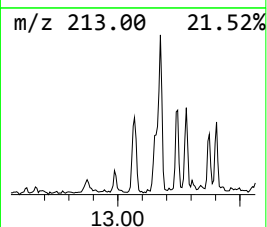
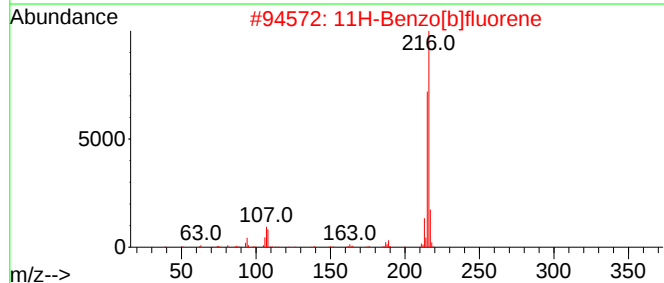
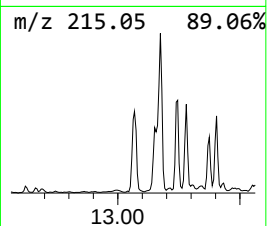
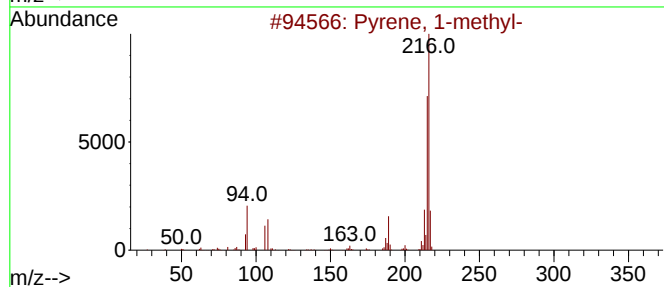
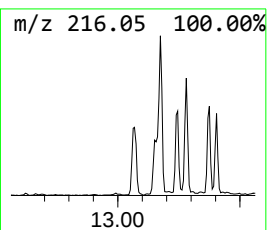
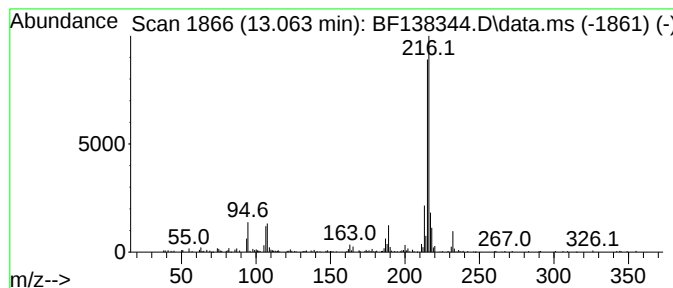
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Pyrene, 1-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD		R.T.	
13.063	2.03 ng	238359	Chrysene-d12		14.045	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Pyrene, 1-methyl-		216	C17H12	002381-21-7	96
2	11H-Benzo[b]fluorene		216	C17H12	000243-17-4	95
3	Fluoranthene, 2-methyl-		216	C17H12	033543-31-6	94
4	Pyrene, 2-methyl-		216	C17H12	003442-78-2	89
5	11H-Benzo[a]fluorene		216	C17H12	000238-84-6	81



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF062524\
Data File : BF138344.D
Acq On : 25 Jun 2024 20:08
Operator : CG\JU
Sample : P3021-10 10X
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
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ClientSampleId :
ARS20-0015

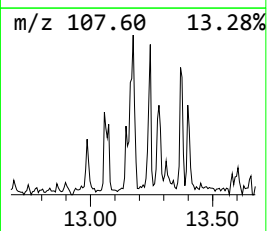
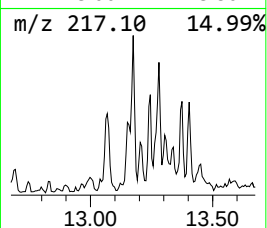
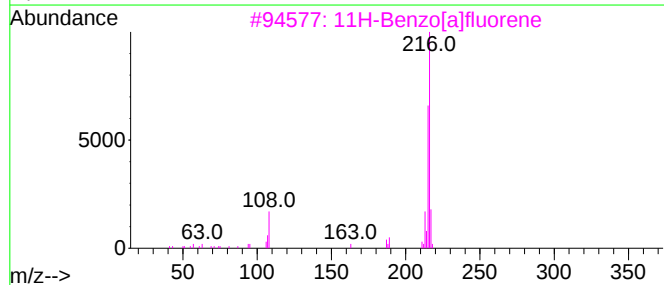
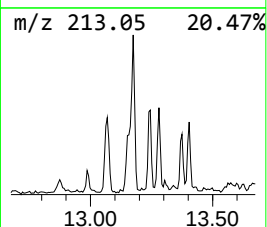
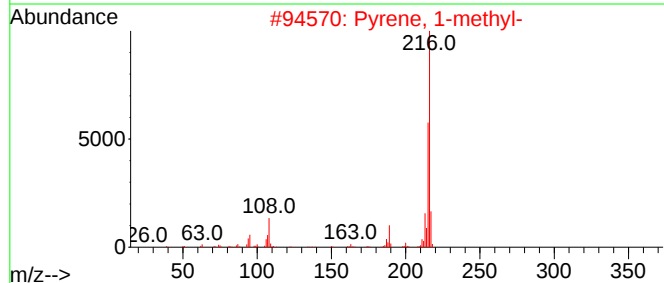
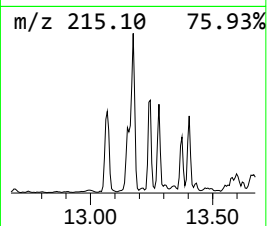
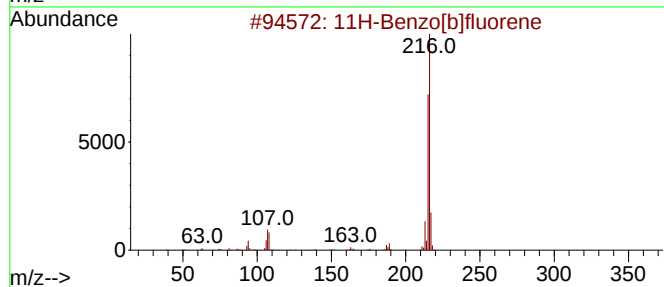
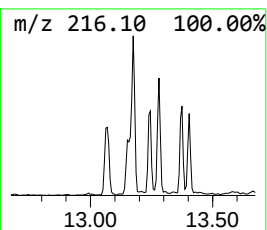
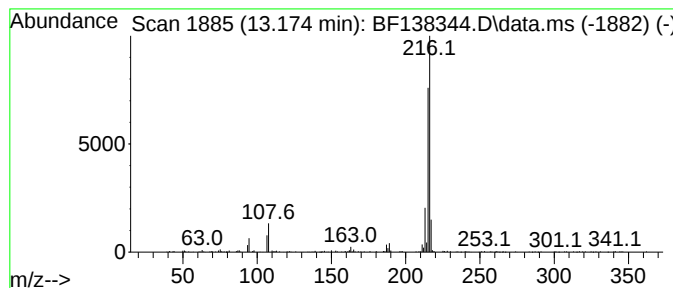
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 11H-Benzo[b]fluorene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.174	2.22 ng	259693	Chrysene-d12	14.045

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			11H-Benzo[b]fluorene	216	C17H12	000243-17-4	94
2			Pyrene, 1-methyl-	216	C17H12	002381-21-7	87
3			11H-Benzo[a]fluorene	216	C17H12	000238-84-6	83
4			Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	83
5			7H-Benzo[c]fluorene	216	C17H12	000205-12-9	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF062524\
Data File : BF138344.D
Acq On : 25 Jun 2024 20:08
Operator : CG\JU
Sample : P3021-10 10X
Misc :
ALS Vial : 21 Sample Multiplier: 1

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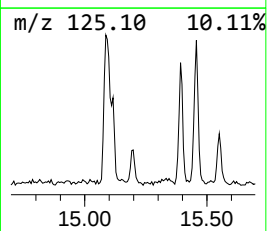
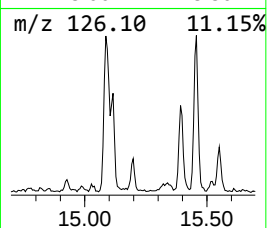
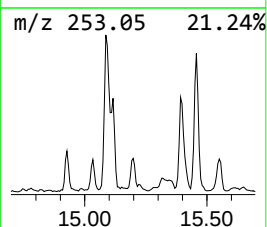
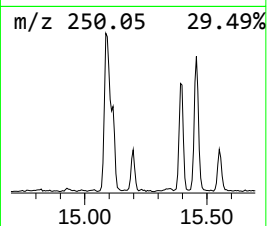
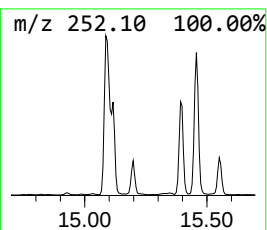
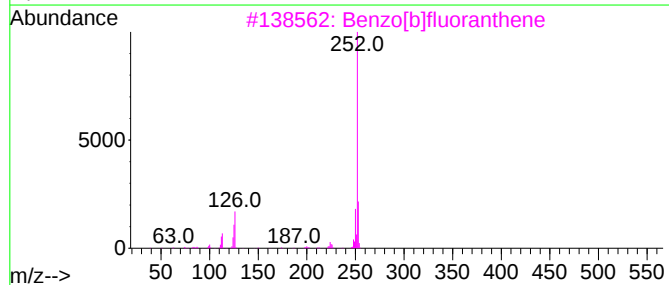
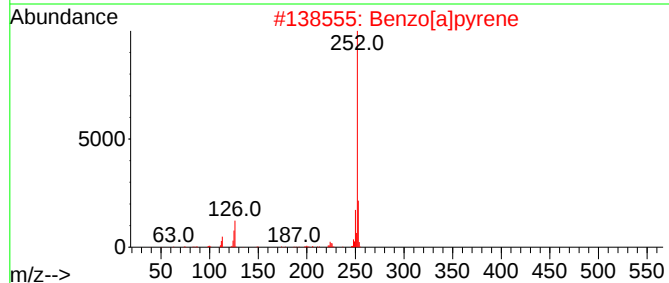
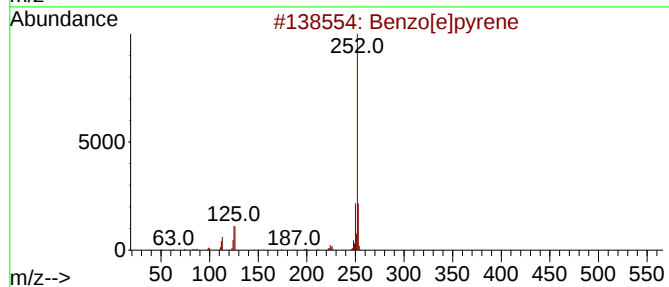
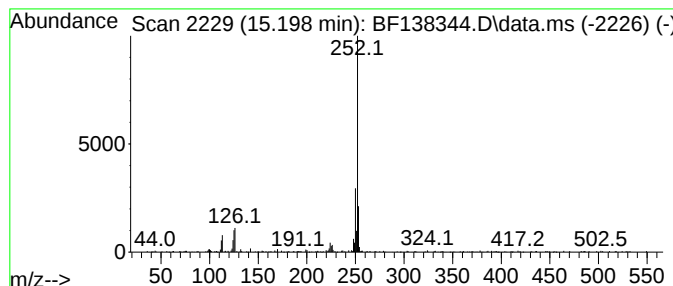
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Benzo[e]pyrene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.198	3.01 ng	124338	Perylene-d12	15.521

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[e]pyrene	252	C20H12	000192-97-2	97
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	94
5			Perylene	252	C20H12	000198-55-0	89



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF062524\
Data File : BF138344.D
Acq On : 25 Jun 2024 20:08
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Misc :
ALS Vial : 21 Sample Multiplier: 1

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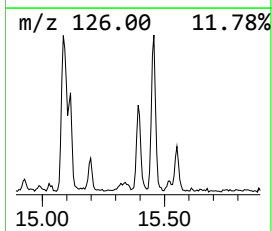
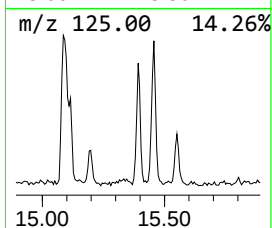
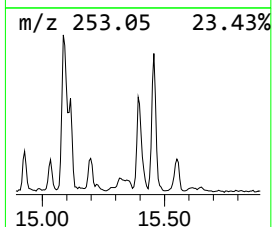
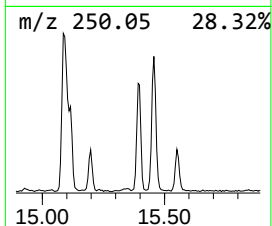
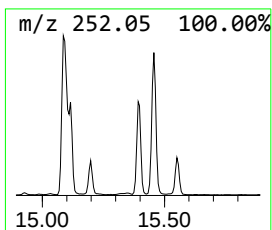
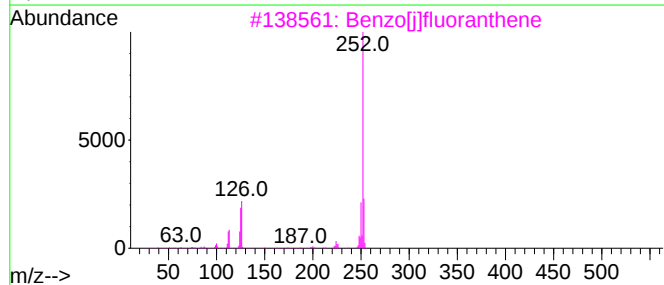
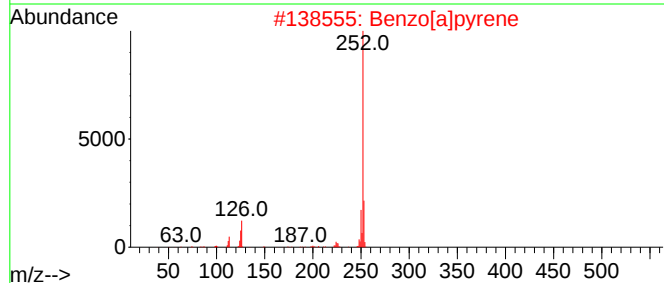
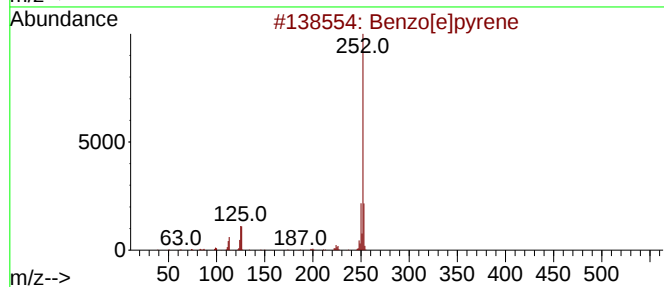
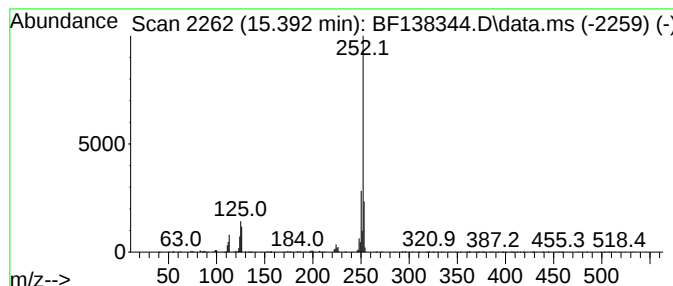
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Benzo[j]fluoranthene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.392	8.86 ng	366020	Perylene-d12	15.521

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzo[e]pyrene	252	C20H12	000192-97-2	98
2		Benzo[a]pyrene	252	C20H12	000050-32-8	98
3		Benzo[j]fluoranthene	252	C20H12	000205-82-3	95
4		Benzo[b]fluoranthene	252	C20H12	000205-99-2	93
5		Perylene	252	C20H12	000198-55-0	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF062524\
Data File : BF138344.D
Acq On : 25 Jun 2024 20:08
Operator : CG\JU
Sample : P3021-10 10X
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
ARS20-0015

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF061224.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-metho...	2.146	8.1	ng	461469	1	6.887	1143710	20.0
Phenanthrene, 2...	11.910	2.9	ng	195945	4	11.404	1331710	20.0
Anthracene, 2-m...	11.933	3.5	ng	234683	4	11.404	1331710	20.0
4H-Cyclopenta[d...	12.016	5.1	ng	338343	4	11.404	1331710	20.0
Phenanthrene, 2...	12.457	2.5	ng	165406	4	11.404	1331710	20.0
Cyclopenta(def)...	12.539	2.0	ng	135686	4	11.404	1331710	20.0
Pyrene, 1-methyl-	13.063	2.0	ng	238359	5	14.045	2343870	20.0
11H-Benzo[b]flu...	13.174	2.2	ng	259693	5	14.045	2343870	20.0
Benzo[e]pyrene	15.198	3.0	ng	124338	6	15.521	826494	20.0
Benzo[j]fluoran...	15.392	8.9	ng	366020	6	15.521	826494	20.0