

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF052422\
 Data File : BF128442.D
 Acq On : 24 May 2022 18:44
 Operator : CG\JU
 Sample : N2981-01
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-13(0.5-1)

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-BF052022.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF128442.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.569	383	388	400	rBV	832349	934955	41.34%	4.395%
2	5.087	473	476	493	rVB	24601	39120	1.73%	0.184%
3	5.475	538	542	562	rBV	1985670	2047935	90.54%	9.626%
4	6.487	709	714	726	rBV	2074665	2139137	94.58%	10.055%
5	6.587	729	731	740	rVB	42081	48905	2.16%	0.230%
6	6.845	771	775	781	rBV	639180	563419	24.91%	2.648%
7	7.410	867	871	881	rBV	1441224	1473433	65.14%	6.926%
8	8.128	989	993	996	rBV	856042	742784	32.84%	3.491%
9	8.839	1111	1114	1120	rVB	24267	23930	1.06%	0.112%
10	9.204	1171	1176	1179	rBV	2490853	2261839	100.00%	10.632%
11	9.269	1183	1187	1190	rVB	38725	32913	1.46%	0.155%
12	9.745	1264	1268	1273	rBV	35605	33083	1.46%	0.156%
13	9.880	1288	1291	1295	rVB	870242	756910	33.46%	3.558%
14	10.298	1355	1362	1365	rBV3	27805	40685	1.80%	0.191%
15	10.433	1381	1385	1391	rVB3	13986	25996	1.15%	0.122%
16	10.675	1423	1426	1440	rBV	1262557	1155491	51.09%	5.431%
17	10.775	1440	1443	1448	rBV2	31042	42876	1.90%	0.202%
18	11.104	1495	1499	1502	rBV4	17040	25815	1.14%	0.121%
19	11.186	1509	1513	1516	rBV3	22673	24889	1.10%	0.117%
20	11.222	1516	1519	1522	rBV2	32716	29177	1.29%	0.137%
21	11.375	1541	1545	1547	rBV	832872	674548	29.82%	3.171%
22	11.398	1547	1549	1554	rVB	205913	163328	7.22%	0.768%
23	11.451	1554	1558	1561	rBV	38511	40891	1.81%	0.192%
24	11.616	1581	1586	1589	rBV3	25858	37973	1.68%	0.178%
25	11.874	1628	1630	1633	rBV	56734	59900	2.65%	0.282%
26	11.904	1633	1635	1639	rVV	91442	81911	3.62%	0.385%
27	11.957	1640	1644	1646	rVV	31983	35957	1.59%	0.169%
28	11.992	1646	1650	1652	rVV2	67281	86161	3.81%	0.405%
29	12.010	1652	1653	1657	rVB	42225	33293	1.47%	0.156%
30	12.169	1677	1680	1685	rVV	39601	40122	1.77%	0.189%
31	12.216	1685	1688	1691	rVB3	27256	29470	1.30%	0.139%
32	12.322	1700	1706	1709	rBV4	20516	29497	1.30%	0.139%
33	12.427	1720	1724	1726	rBV	66467	65454	2.89%	0.308%
34	12.521	1736	1740	1743	rBV2	31526	43918	1.94%	0.206%
35	12.592	1748	1752	1755	rBV	470699	395438	17.48%	1.859%
36	12.692	1766	1769	1771	rBV	27434	27558	1.22%	0.130%

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

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 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

37	12.804	1785	1788	1789	rBV	85601	81009	3.58%	0.381%
38	12.821	1789	1791	1797	rVV	402297	363910	16.09%	1.711%
39	12.874	1797	1800	1804	rVB2	44655	48953	2.16%	0.230%
40	12.957	1806	1814	1818	rBV	1697768	1669943	73.83%	7.850%
41	13.039	1823	1828	1833	rBV3	52905	97234	4.30%	0.457%
42	13.151	1839	1847	1849	rBV2	90707	135999	6.01%	0.639%
43	13.216	1855	1858	1861	rBV	48114	42287	1.87%	0.199%
44	13.257	1861	1865	1868	rVV2	58467	72061	3.19%	0.339%
45	13.351	1878	1881	1883	rBV	40577	35013	1.55%	0.165%
46	13.380	1883	1886	1889	rBV	36232	35987	1.59%	0.169%
47	13.516	1906	1909	1912	rBV4	28452	28469	1.26%	0.134%
48	13.574	1917	1919	1925	rVB3	26939	35618	1.57%	0.167%
49	13.668	1931	1935	1938	rBV	72072	69413	3.07%	0.326%
50	13.774	1951	1953	1955	rBV	62403	45581	2.02%	0.214%
51	13.798	1955	1957	1959	rVB	48216	33713	1.49%	0.158%
52	13.827	1959	1962	1963	rBV	43354	37189	1.64%	0.175%
53	13.845	1963	1965	1968	rVB2	47828	42697	1.89%	0.201%
54	13.874	1968	1970	1977	rBV	78016	115640	5.11%	0.544%
55	14.021	1991	1995	1998	rBV2	698158	925995	40.94%	4.353%
56	14.051	1998	2000	2007	rVB	334731	285786	12.64%	1.343%
57	14.110	2007	2010	2012	rBV3	43027	50744	2.24%	0.239%
58	14.163	2016	2019	2020	rBV2	32623	31723	1.40%	0.149%
59	14.180	2020	2022	2027	rVB4	45760	46346	2.05%	0.218%
60	14.374	2053	2055	2057	rBV	28747	27226	1.20%	0.128%
61	14.410	2057	2061	2065	rVB	87472	106725	4.72%	0.502%
62	14.445	2065	2067	2069	rBV	49708	34628	1.53%	0.163%
63	14.539	2080	2083	2085	rBV	47190	45488	2.01%	0.214%
64	14.915	2144	2147	2150	rBV2	56356	64602	2.86%	0.304%
65	15.074	2170	2174	2177	rBV	347750	491987	21.75%	2.313%
66	15.186	2189	2193	2197	rBV	95344	116566	5.15%	0.548%
67	15.268	2205	2207	2213	rBV4	22782	41914	1.85%	0.197%
68	15.380	2222	2226	2233	rVB	230778	271668	12.01%	1.277%
69	15.445	2233	2237	2242	rVV	286728	341974	15.12%	1.607%
70	15.510	2243	2248	2251	rVV	337639	443128	19.59%	2.083%
71	15.539	2251	2253	2260	rVB2	89377	133697	5.91%	0.628%
72	17.009	2497	2503	2510	rVB	139628	285730	12.63%	1.343%
73	17.186	2531	2533	2539	rVB2	21815	32221	1.42%	0.151%
74	17.462	2576	2580	2594	rVB	98868	216525	9.57%	1.018%

Sum of corrected areas: 21274100

Instrument :

BNA_F

ClientSampleId :

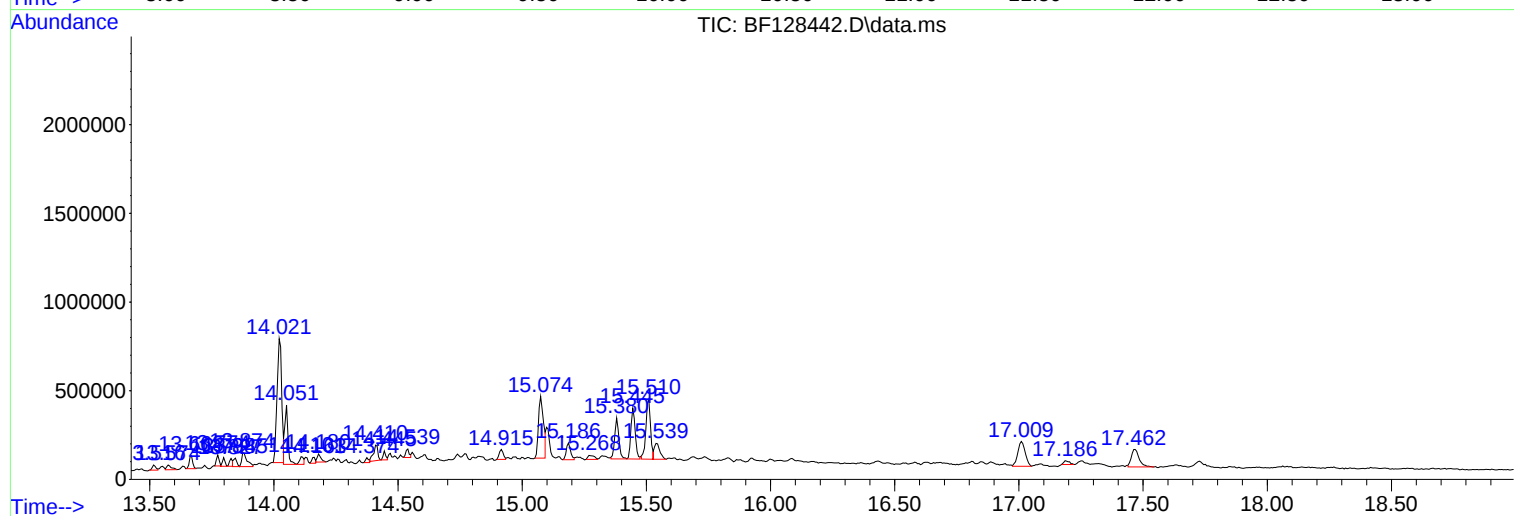
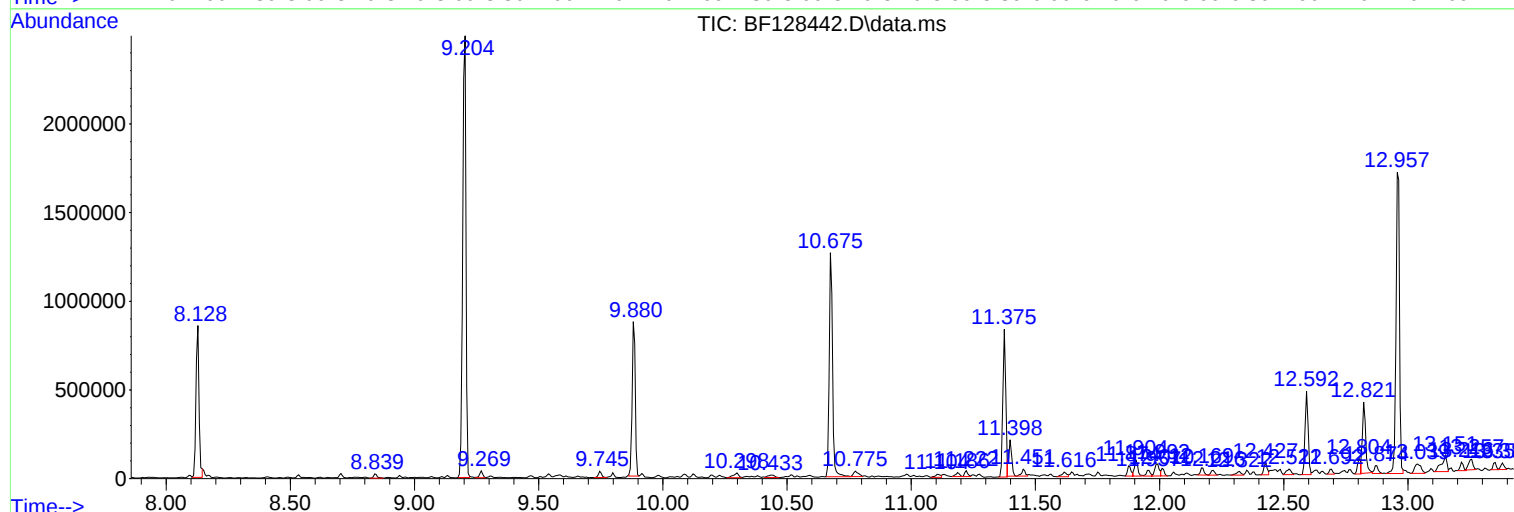
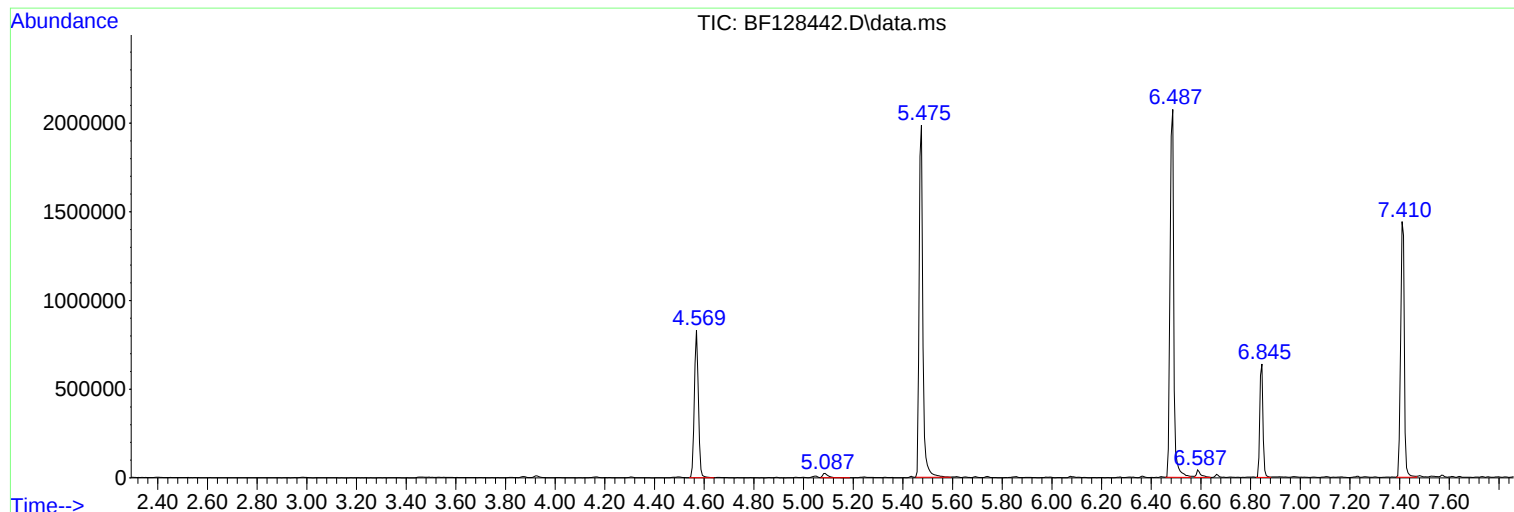
SB-13(0.5-1)

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Instrument :
BNA_F
ClientSampleId :
SB-13(0.5-1)

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF052022.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF052422\
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Operator : CG\JU
Sample : N2981-01
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ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_F
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SB-13(0.5-1)

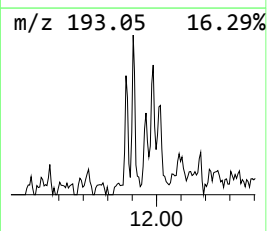
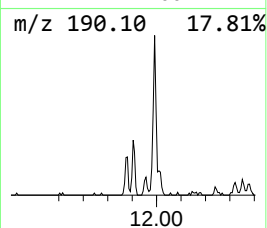
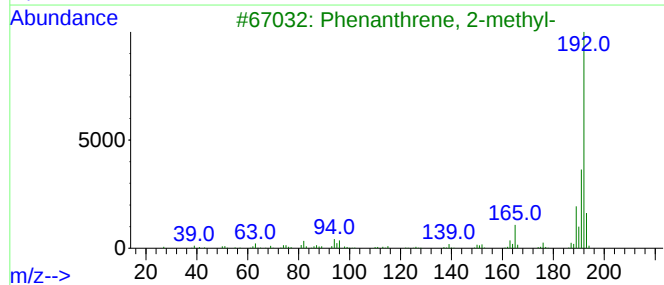
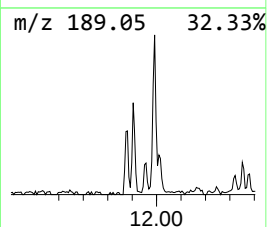
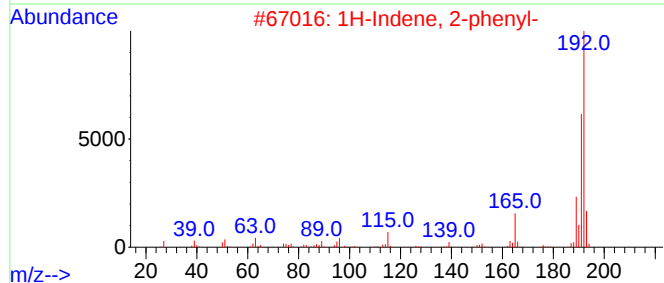
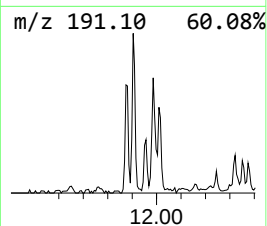
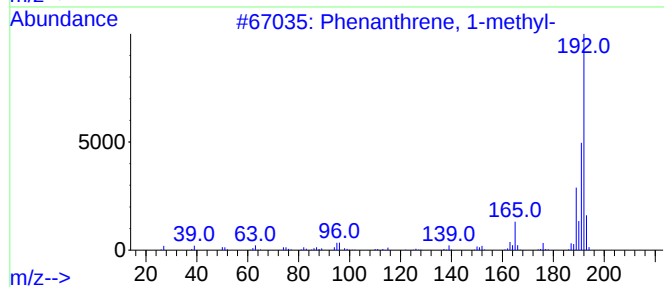
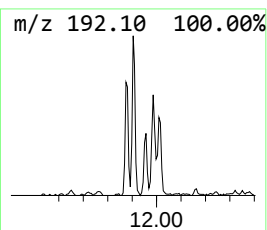
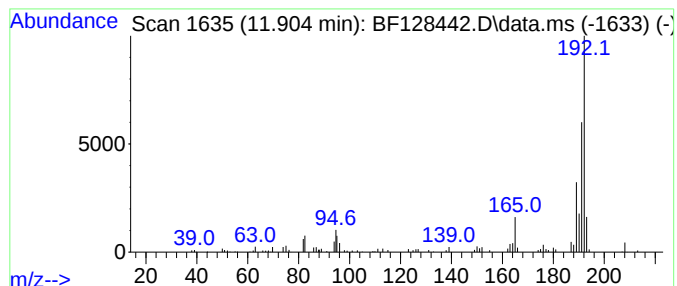
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 Phenanthrene, 1-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.904	2.43 ng	81911	Phenanthrene-d10	11.374

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	97
2			1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	94
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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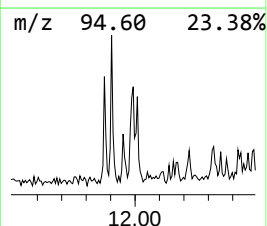
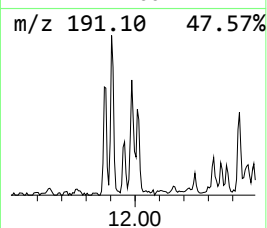
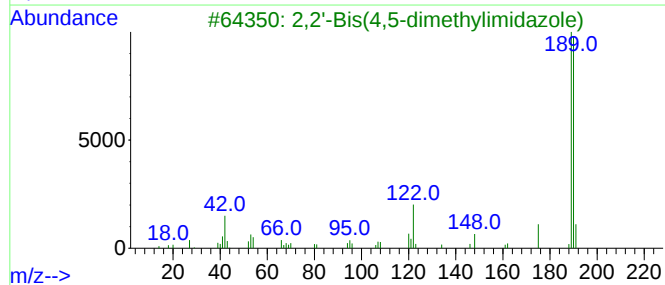
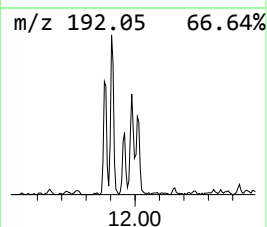
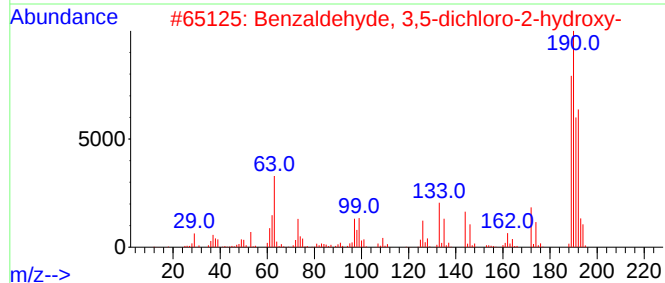
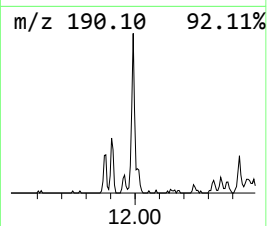
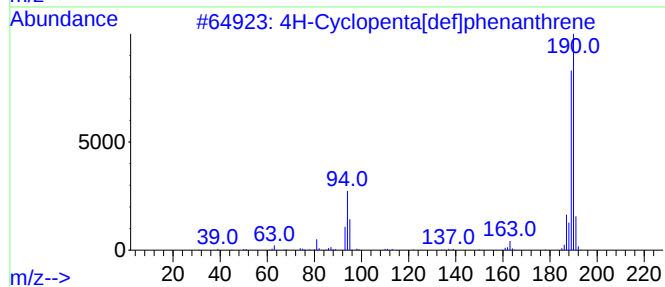
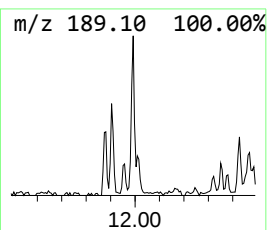
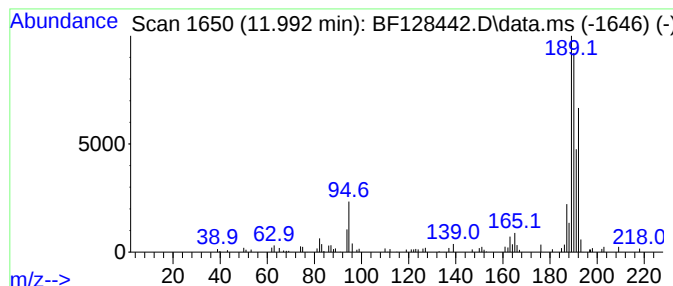
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF052022.M
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 4H-Cyclopenta[def]phenanthrene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.992	2.55 ng	86161	Phenanthrene-d10	11.374

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	76
2			Benzaldehyde, 3,5-dichloro-2-hyd...	190	C7H4Cl2O2	000090-60-8	50
3			2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	43
4			6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	42
5			4,6-Dichloro-2-ethylpyrimidin-5-...	191	C6H7Cl2N3	006237-96-3	38



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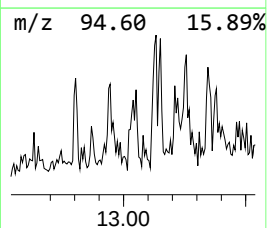
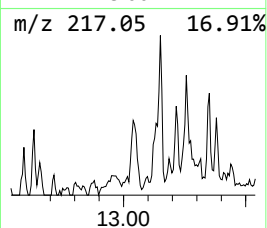
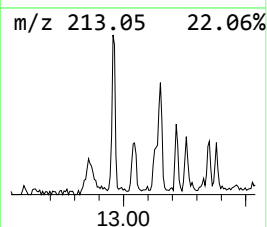
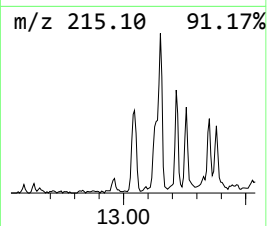
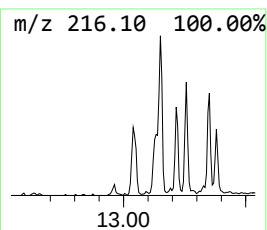
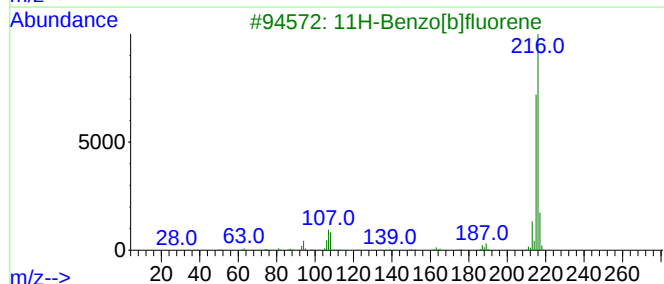
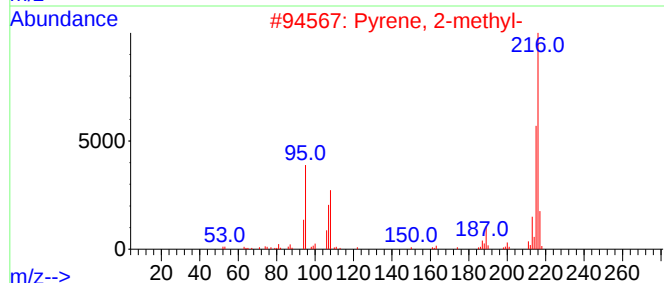
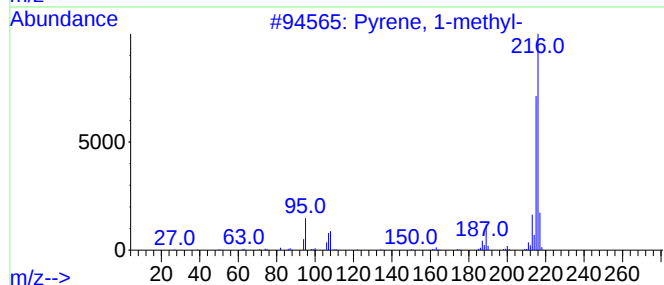
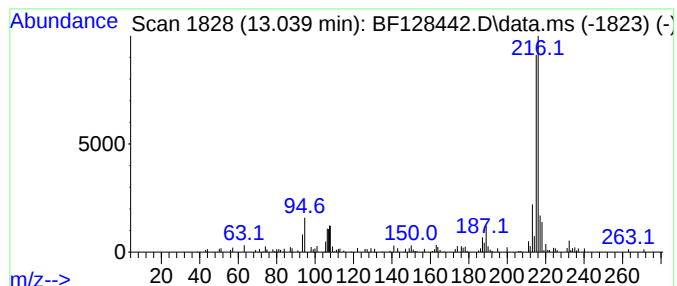
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Pyrene, 1-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.039	2.10 ng	97234	Chrysene-d12	14.027

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pyrene, 1-methyl-	216	C17H12	002381-21-7	96
2			Pyrene, 2-methyl-	216	C17H12	003442-78-2	93
3			11H-Benzo[b]fluorene	216	C17H12	000243-17-4	93
4			Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	89
5			11H-Benzo[a]fluorene	216	C17H12	000238-84-6	76



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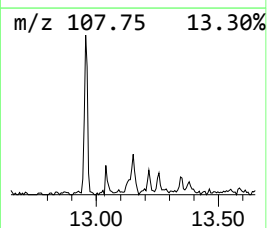
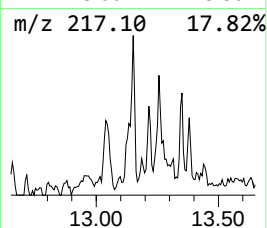
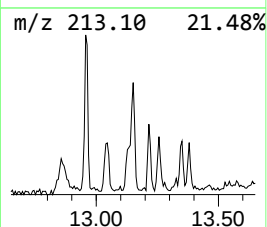
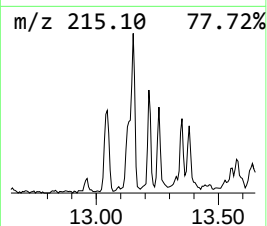
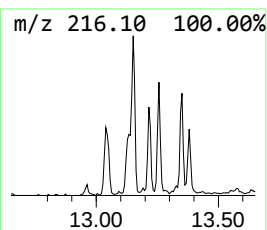
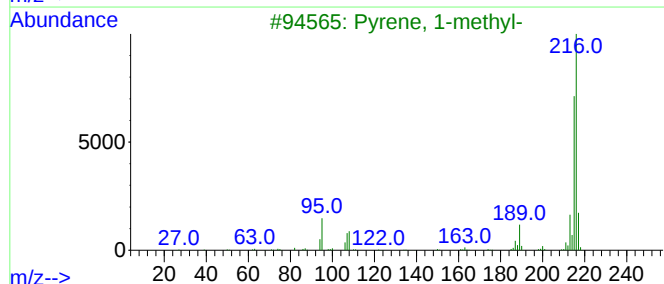
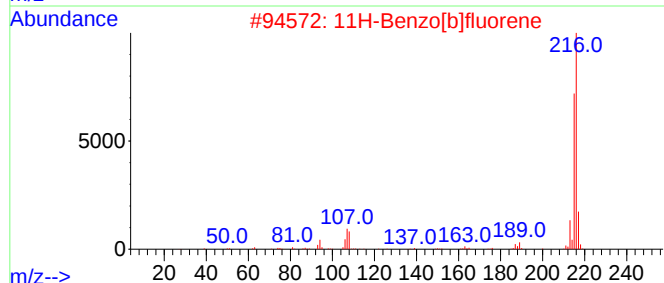
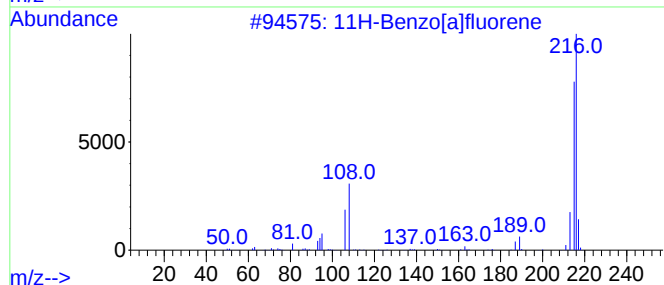
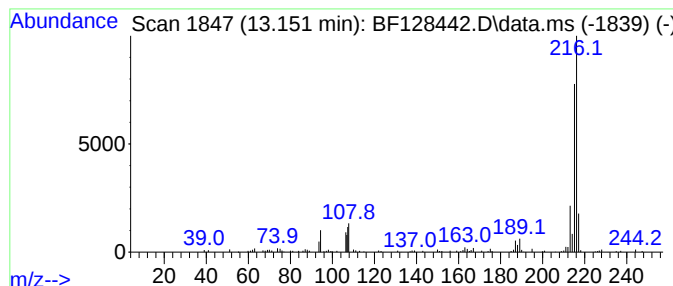
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 11H-Benzo[a]fluorene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.151	2.94 ng	135999	Chrysene-d12	14.027

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			11H-Benzo[a]fluorene	216	C17H12	000238-84-6	94
2			11H-Benzo[b]fluorene	216	C17H12	000243-17-4	94
3			Pyrene, 1-methyl-	216	C17H12	002381-21-7	93
4			Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	90
5			7H-Benzo[c]fluorene	216	C17H12	000205-12-9	74



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF052422\
Data File : BF128442.D
Acq On : 24 May 2022 18:44
Operator : CG\JU
Sample : N2981-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-13(0.5-1)

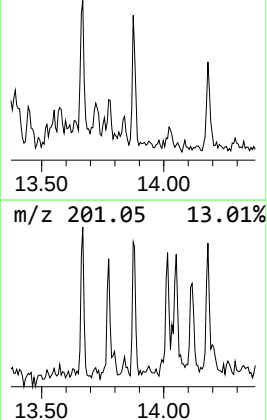
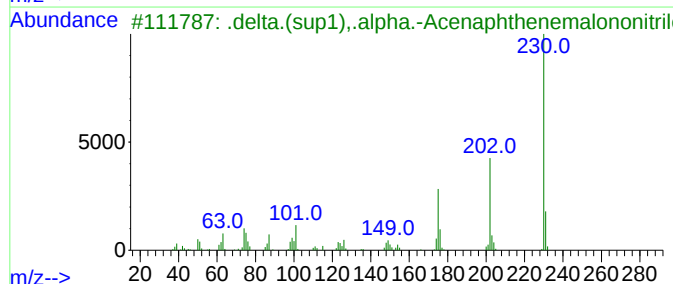
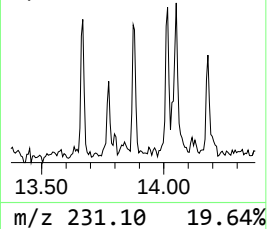
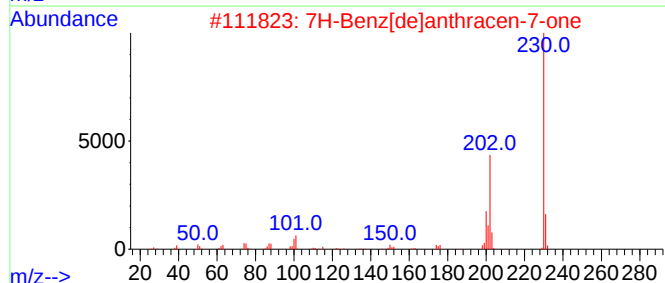
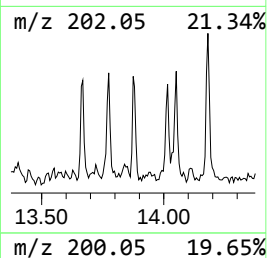
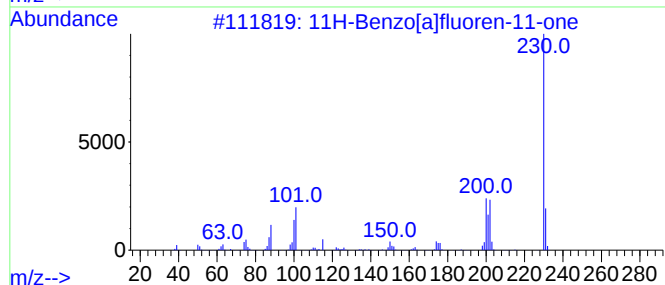
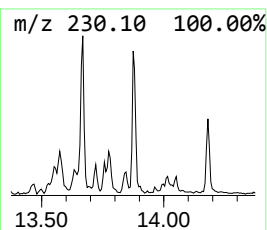
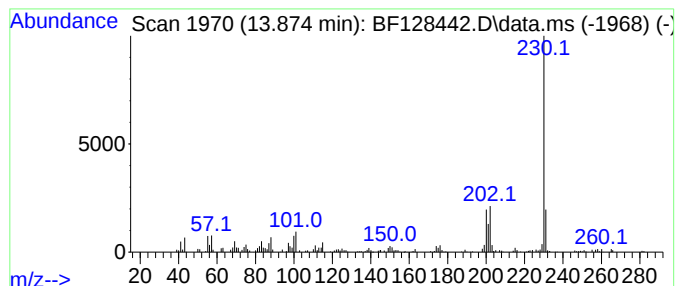
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 11H-Benzo[a]fluoren-11-one Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.874	2.50 ng	115640	Chrysene-d12	14.027

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			11H-Benzo[a]fluoren-11-one	230	C17H10O	000479-79-8	93
2			7H-Benz[de]anthracen-7-one	230	C17H10O	000082-05-3	90
3			.delta.(sup1),.alpha.-Acenaphthe...	230	C15H6N2O	014619-86-4	59
4			1-Pyrenecarboxaldehyde	230	C17H10O	003029-19-4	56
5			2-Chloro-9-xanthenone	230	C13H7ClO2	013210-15-6	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF052422\
Data File : BF128442.D
Acq On : 24 May 2022 18:44
Operator : CG\JU
Sample : N2981-01
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ALS Vial : 15 Sample Multiplier: 1

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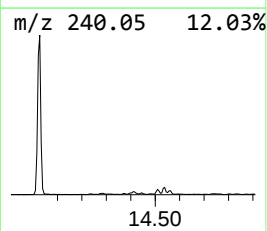
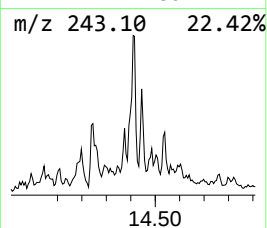
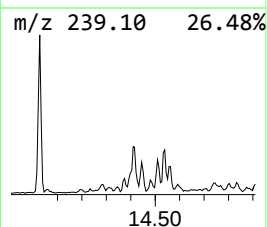
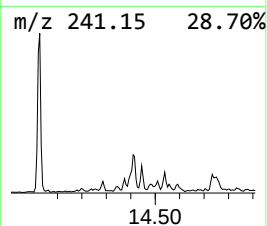
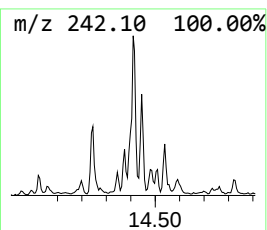
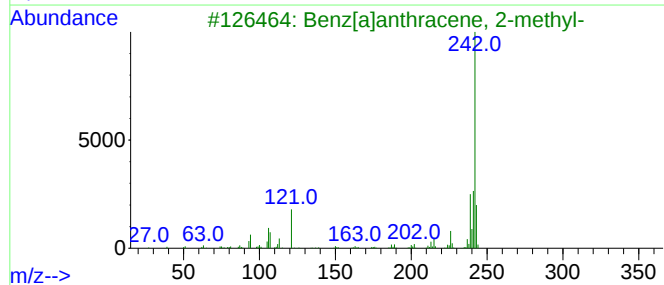
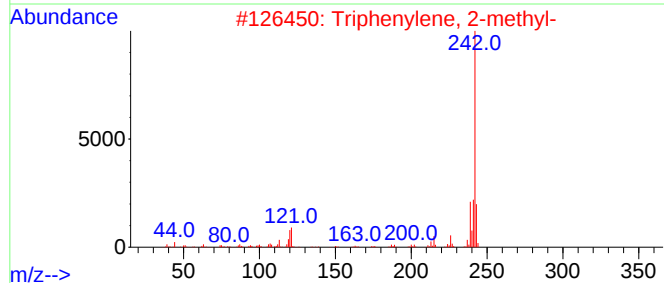
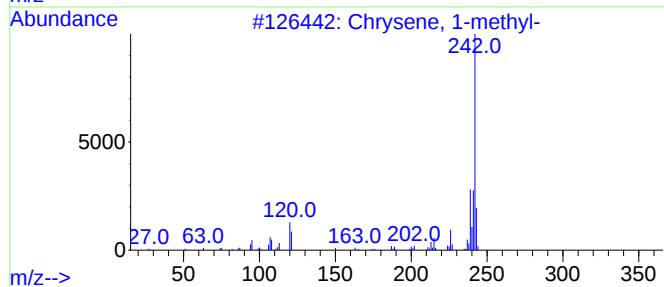
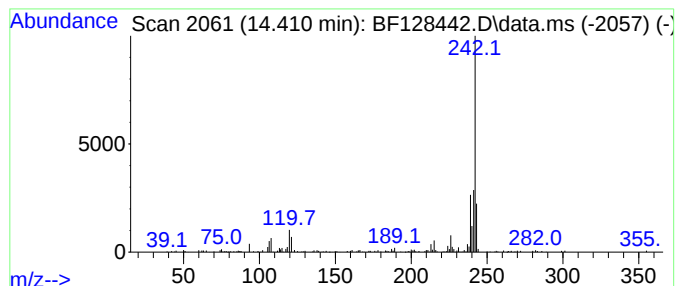
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Chrysene, 1-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.410	2.31 ng	106725	Chrysene-d12	14.027

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Chrysene, 1-methyl-	242	C19H14	003351-28-8	96
2			Triphenylene, 2-methyl-	242	C19H14	001705-84-6	95
3			Benz[a]anthracene, 2-methyl-	242	C19H14	002498-76-2	93
4			Chrysene, 6-methyl-	242	C19H14	001705-85-7	91
5			Chrysene, 5-methyl-	242	C19H14	003697-24-3	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF052422\
Data File : BF128442.D
Acq On : 24 May 2022 18:44
Operator : CG\JU
Sample : N2981-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

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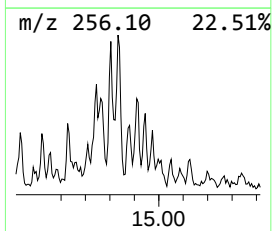
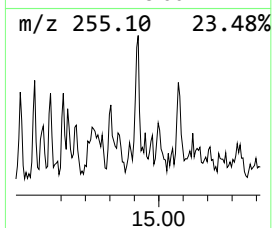
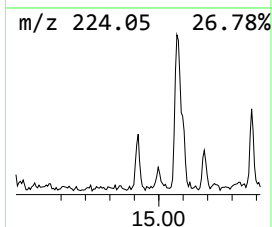
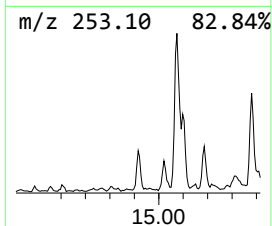
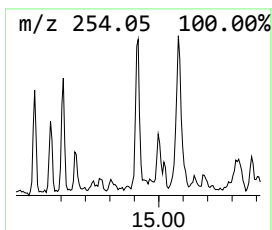
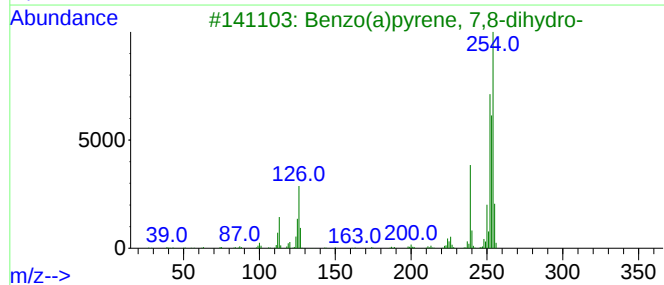
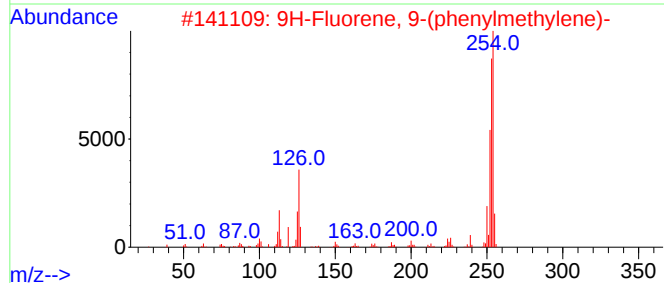
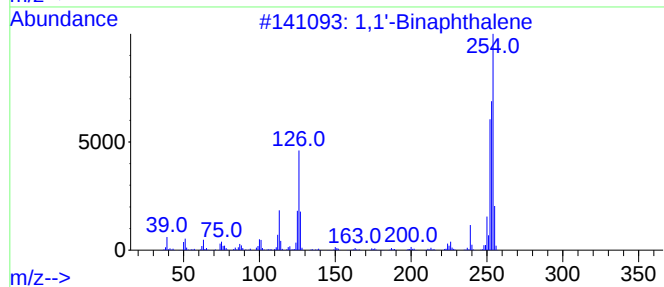
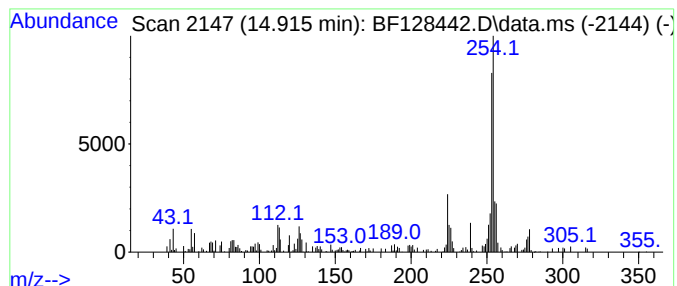
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 1,1'-Binaphthalene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.916	2.92 ng	64602	Perylene-d12	15.509

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,1'-Binaphthalene	254	C20H14	000604-53-5	76
2			9H-Fluorene, 9-(phenylmethylene)-	254	C20H14	001836-87-9	70
3			Benzo(a)pyrene, 7,8-dihydro-	254	C20H14	017573-23-8	68
4			1,2'-Binaphthalene	254	C20H14	004325-74-0	68
5			Benzo[a]pyrene, 4,5-dihydro-	254	C20H14	057652-66-1	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF052422\
Data File : BF128442.D
Acq On : 24 May 2022 18:44
Operator : CG\JU
Sample : N2981-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

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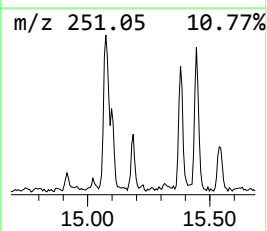
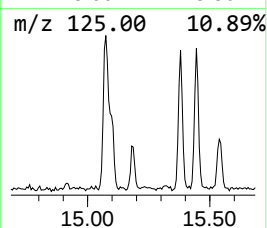
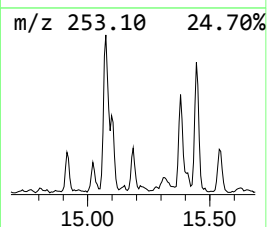
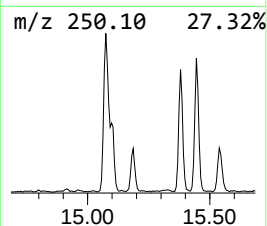
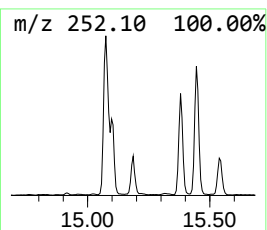
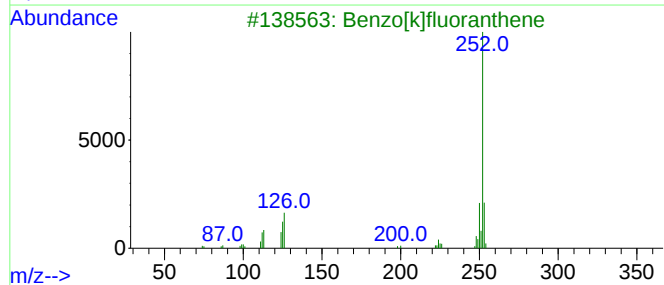
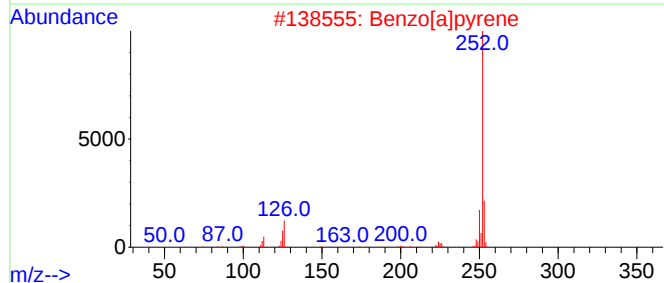
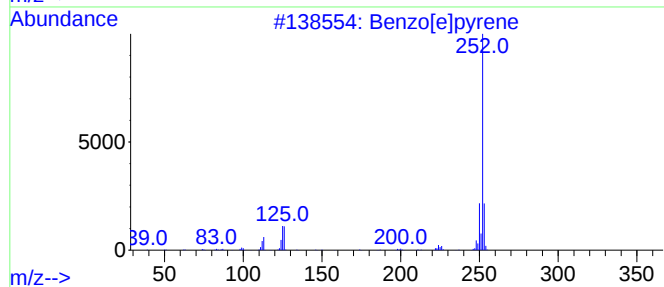
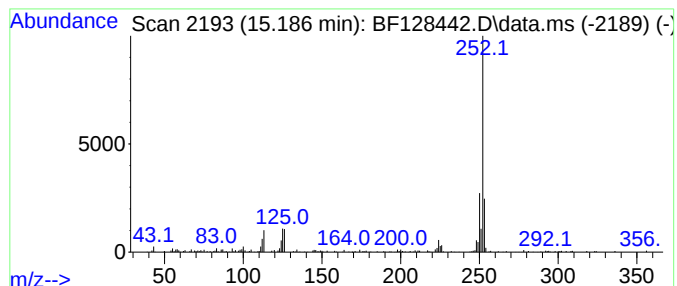
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Benzo[e]pyrene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.186	5.26 ng	116566	Perylene-d12	15.509

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	95
3			Benzo[k]fluoranthene	252	C20H12	000207-08-9	95
4			Benzo[b]fluoranthene	252	C20H12	000205-99-2	94
5			Perylene	252	C20H12	000198-55-0	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF052422\
Data File : BF128442.D
Acq On : 24 May 2022 18:44
Operator : CG\JU
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Misc :
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Instrument :
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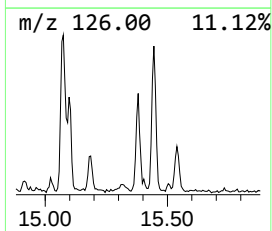
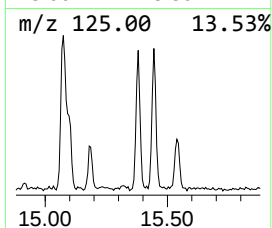
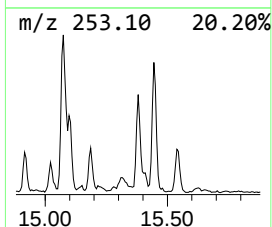
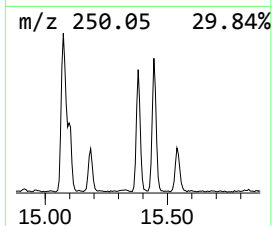
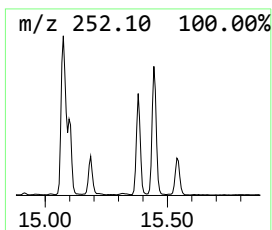
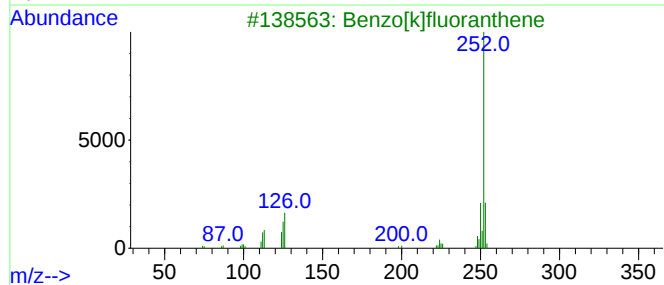
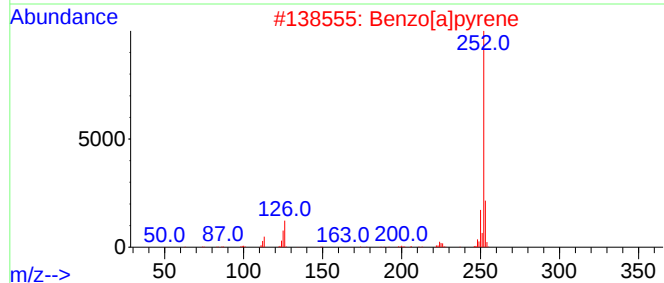
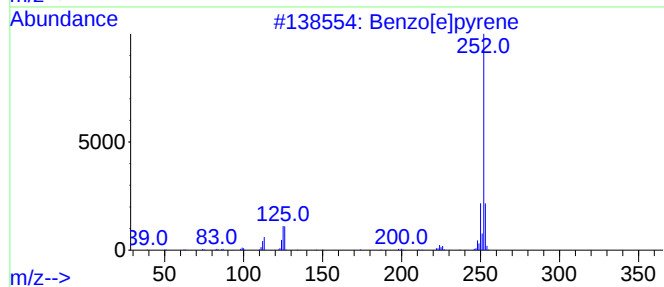
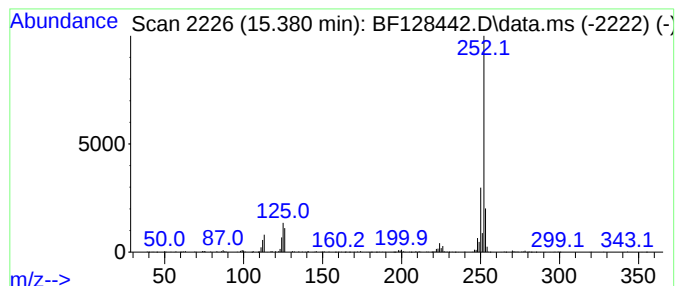
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Benzo[j]fluoranthene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.380	12.26 ng	271668	Perylene-d12	15.509

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	95
3			Benzo[k]fluoranthene	252	C20H12	000207-08-9	94
4			Benzo[b]fluoranthene	252	C20H12	000205-99-2	93
5			Benzo[j]fluoranthene	252	C20H12	000205-82-3	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF052422\
Data File : BF128442.D
Acq On : 24 May 2022 18:44
Operator : CG\JU
Sample : N2981-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-13(0.5-1)

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF052022.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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					#	RT	Resp	Conc
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4H-Cyclopenta[d...	11.992	2.5	ng	86161	4	11.374	674548	20.0
Pyrene, 1-methyl-	13.039	2.1	ng	97234	5	14.027	925995	20.0
11H-Benzo[a]flu...	13.151	2.9	ng	135999	5	14.027	925995	20.0
11H-Benzo[a]flu...	13.874	2.5	ng	115640	5	14.027	925995	20.0
Chrysene, 1-met...	14.410	2.3	ng	106725	5	14.027	925995	20.0
1,1'-Binaphthalene	14.916	2.9	ng	64602	6	15.509	443128	20.0
Benzo[e]pyrene	15.186	5.3	ng	116566	6	15.509	443128	20.0
Benzo[j]fluoran...	15.380	12.3	ng	271668	6	15.509	443128	20.0