

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF062122\
 Data File : BF128916.D
 Acq On : 22 Jun 2022 00:19
 Operator : CG\JU
 Sample : N3421-04
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-3-0-2-20220620

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-BF060222.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF128916.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.969	451	456	465	rVB	35475	48582	2.15%	0.194%
2	5.392	524	528	540	rBV	2371053	2191797	97.01%	8.770%
3	6.410	697	701	714	rBV	2404317	2259300	100.00%	9.040%
4	6.498	714	716	725	rVB	34274	31167	1.38%	0.125%
5	6.751	756	759	763	rBV	738980	587364	26.00%	2.350%
6	7.322	851	856	859	rBV	1677265	1514353	67.03%	6.059%
7	8.033	972	977	980	rBV	788687	762241	33.74%	3.050%
8	8.057	980	981	986	rVB	178643	102455	4.53%	0.410%
9	8.751	1095	1099	1102	rBV	38764	33067	1.46%	0.132%
10	9.110	1155	1160	1164	rBV	2410182	2243453	99.30%	8.977%
11	9.786	1272	1275	1278	rBV	807998	702522	31.09%	2.811%
12	9.822	1278	1281	1284	rVB	198441	154728	6.85%	0.619%
13	9.992	1307	1310	1315	rBV	79276	63745	2.82%	0.255%
14	10.333	1365	1368	1374	rVB	127971	109679	4.85%	0.439%
15	10.404	1374	1380	1381	rBV	22193	24196	1.07%	0.097%
16	10.580	1406	1410	1422	rBV	1320395	1148021	50.81%	4.594%
17	11.086	1492	1496	1499	rVB	31437	28338	1.25%	0.113%
18	11.169	1508	1510	1517	rVB	43327	39346	1.74%	0.157%
19	11.274	1524	1528	1530	rBV	746040	629211	27.85%	2.518%
20	11.298	1530	1532	1536	rVV	899655	726885	32.17%	2.908%
21	11.351	1536	1541	1544	rVB	244189	214700	9.50%	0.859%
22	11.510	1563	1568	1572	rBV2	126832	124365	5.50%	0.498%
23	11.669	1589	1595	1597	rBV4	28430	27624	1.22%	0.111%
24	11.774	1610	1613	1616	rBV	68065	64602	2.86%	0.258%
25	11.804	1616	1618	1622	rVV2	102637	95320	4.22%	0.381%
26	11.851	1622	1626	1629	rVV2	39261	42101	1.86%	0.168%
27	11.892	1629	1633	1635	rVV	151867	153763	6.81%	0.615%
28	11.910	1635	1636	1642	rVB	42859	30579	1.35%	0.122%
29	12.074	1661	1664	1666	rBV	55915	47279	2.09%	0.189%
30	12.104	1666	1669	1672	rVB	53732	43672	1.93%	0.175%
31	12.327	1703	1707	1710	rBV2	27277	35115	1.55%	0.141%
32	12.416	1719	1722	1726	rBV2	32077	34565	1.53%	0.138%
33	12.492	1731	1735	1738	rVV	1376058	1133828	50.18%	4.537%
34	12.639	1754	1760	1763	rBV2	44196	56797	2.51%	0.227%
35	12.721	1768	1774	1777	rVV	1115029	1172883	51.91%	4.693%
36	12.768	1779	1782	1785	rVB	42166	41129	1.82%	0.165%

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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-BF060222.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

37	12.839	1790	1794	1795	rBV	76191	72748	3.22%	0.291%
38	12.863	1795	1798	1801	rVV	1875332	1547551	68.50%	6.192%
39	12.939	1805	1811	1814	rBV2	58801	102836	4.55%	0.411%
40	12.974	1814	1817	1819	rBV	38684	28717	1.27%	0.115%
41	13.021	1821	1825	1827	rBV4	62287	75575	3.35%	0.302%
42	13.045	1827	1829	1833	rVB	173358	165748	7.34%	0.663%
43	13.086	1833	1836	1838	rBV2	17867	25520	1.13%	0.102%
44	13.116	1838	1841	1843	rBV	160547	131027	5.80%	0.524%
45	13.151	1843	1847	1849	rBV3	96677	105842	4.68%	0.424%
46	13.174	1849	1851	1854	rVB	60619	51567	2.28%	0.206%
47	13.204	1854	1856	1859	rVB	29716	28013	1.24%	0.112%
48	13.245	1859	1863	1865	rBV	46095	42035	1.86%	0.168%
49	13.274	1865	1868	1872	rBV2	53024	57559	2.55%	0.230%
50	13.433	1891	1895	1897	rBV4	36506	46238	2.05%	0.185%
51	13.474	1899	1902	1904	rVB2	37424	42114	1.86%	0.169%
52	13.533	1909	1912	1914	rBV2	29239	31780	1.41%	0.127%
53	13.557	1914	1916	1920	rVB	63970	65287	2.89%	0.261%
54	13.668	1931	1935	1937	rBV	126687	119257	5.28%	0.477%
55	13.692	1937	1939	1942	rVV	99818	95958	4.25%	0.384%
56	13.721	1942	1944	1945	rVV	87103	77297	3.42%	0.309%
57	13.768	1949	1952	1957	rVB2	82382	81655	3.61%	0.327%
58	13.839	1959	1964	1969	rBV	42087	77757	3.44%	0.311%
59	13.915	1969	1977	1980	rVV2	892548	1394430	61.72%	5.579%
60	13.945	1980	1982	1989	rVV	599877	521358	23.08%	2.086%
61	14.021	1989	1995	1998	rVV2	144711	198164	8.77%	0.793%
62	14.080	2001	2005	2008	rVV5	40797	77276	3.42%	0.309%
63	14.115	2008	2011	2013	rVV	71945	72327	3.20%	0.289%
64	14.151	2015	2017	2020	rVB	76545	61488	2.72%	0.246%
65	14.268	2034	2037	2039	rBV2	43622	40234	1.78%	0.161%
66	14.304	2039	2043	2046	rVV	94783	104466	4.62%	0.418%
67	14.339	2046	2049	2053	rVB2	48556	43752	1.94%	0.175%
68	14.386	2053	2057	2058	rBV4	45383	48405	2.14%	0.194%
69	14.404	2058	2060	2062	rVV	44514	38428	1.70%	0.154%
70	14.433	2062	2065	2066	rVV	46159	45924	2.03%	0.184%
71	14.451	2066	2068	2071	rVB2	72204	58627	2.59%	0.235%
72	14.615	2093	2096	2098	rBV2	48419	43880	1.94%	0.176%
73	14.798	2123	2127	2133	rVB3	37434	65506	2.90%	0.262%
74	14.951	2148	2153	2156	rBV	454670	683822	30.27%	2.736%
75	15.051	2167	2170	2174	rBV	67213	68639	3.04%	0.275%
76	15.139	2182	2185	2186	rBV2	28381	30441	1.35%	0.122%

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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

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Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

77	15.239	2198	2202	2209	rVB	239444	302901	13.41%	1.212%
78	15.304	2209	2213	2218	rVV	314191	379924	16.82%	1.520%
79	15.362	2218	2223	2226	rVV	326904	404355	17.90%	1.618%
80	15.392	2226	2228	2234	rVB	102263	130952	5.80%	0.524%
81	16.780	2459	2464	2472	rVB	122927	213343	9.44%	0.854%
82	17.203	2532	2536	2547	rVB	79603	172569	7.64%	0.690%

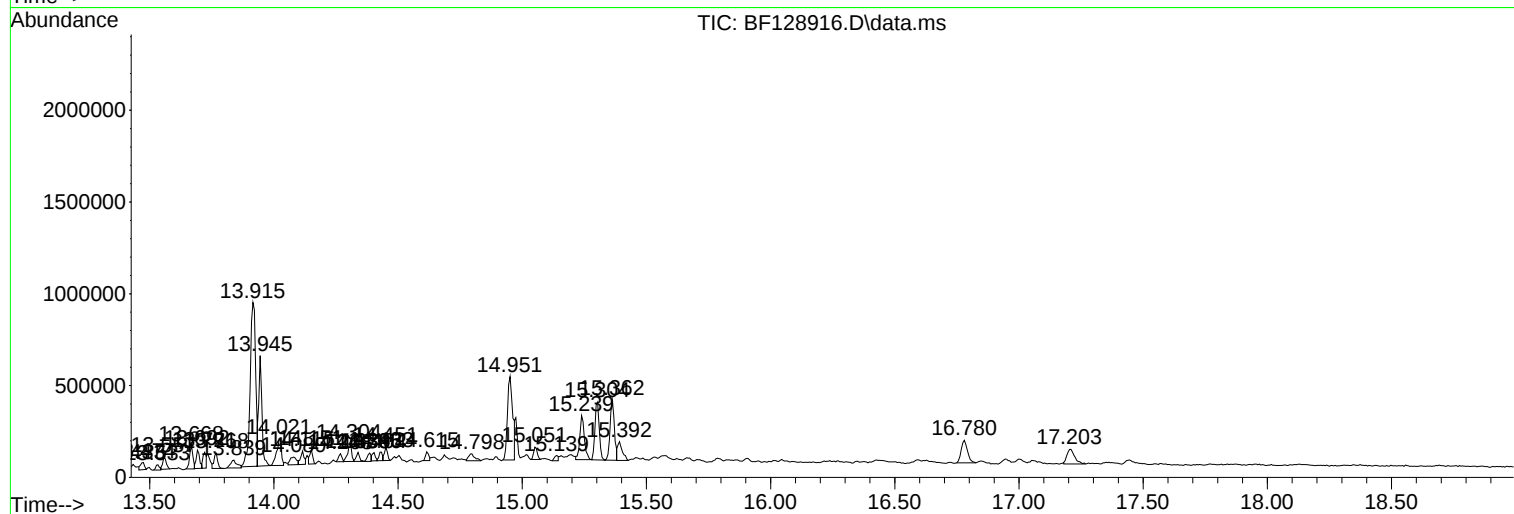
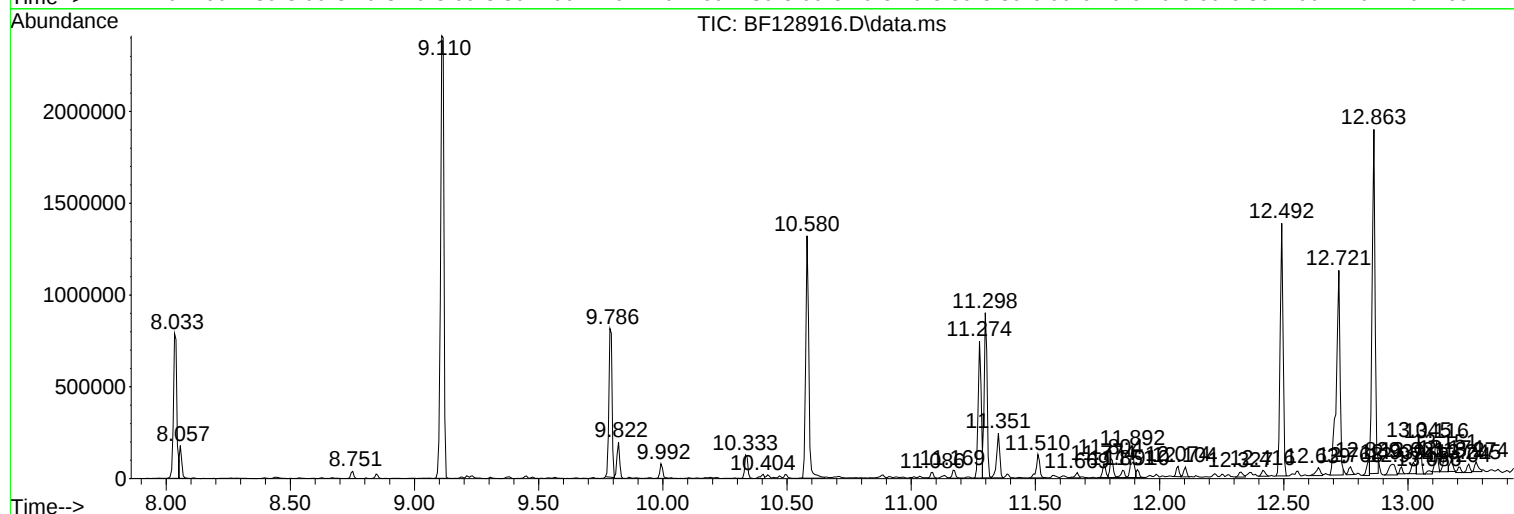
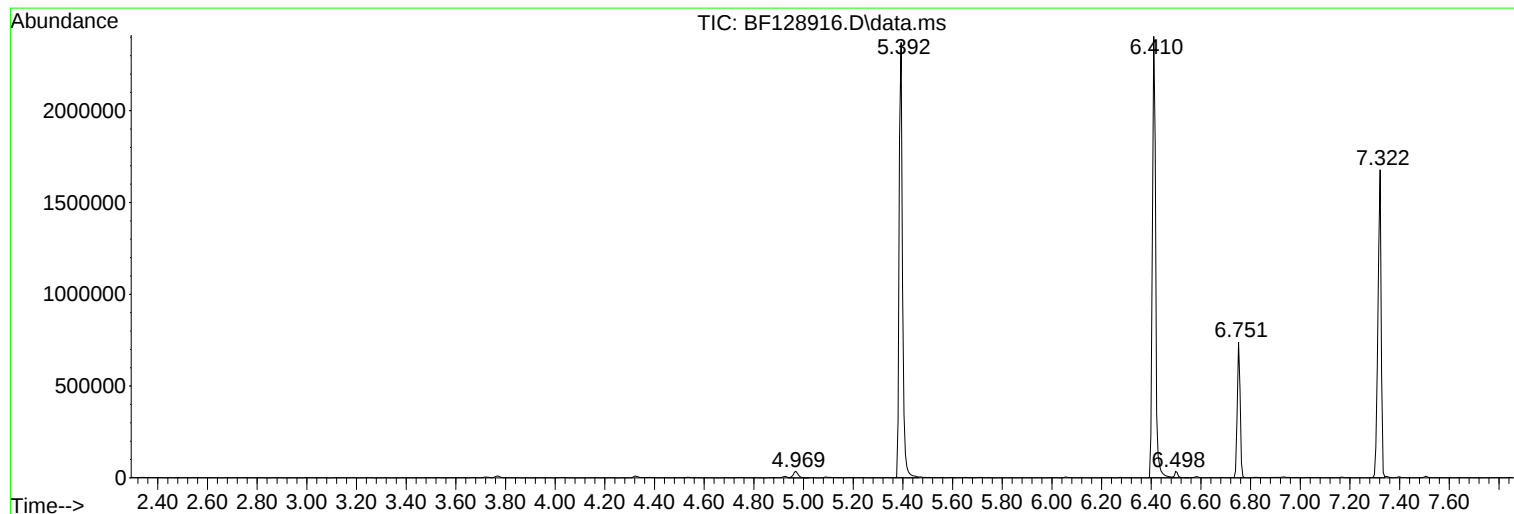
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF060222.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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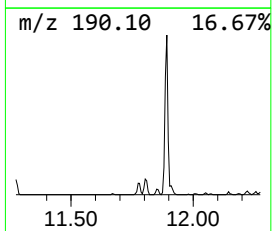
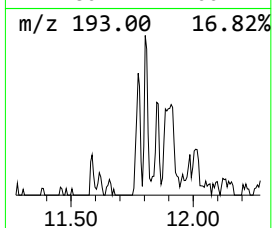
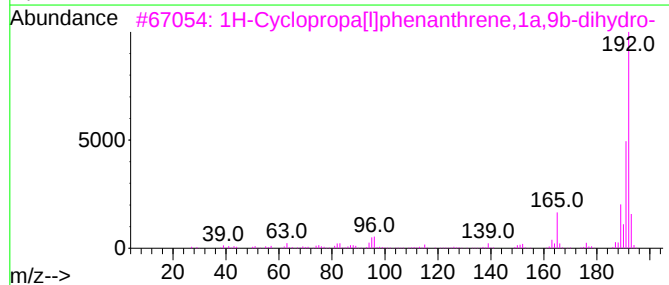
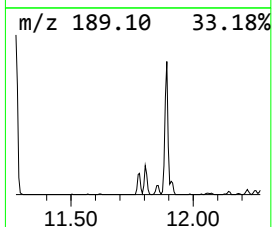
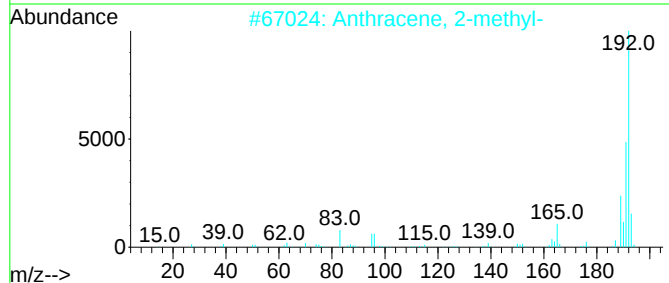
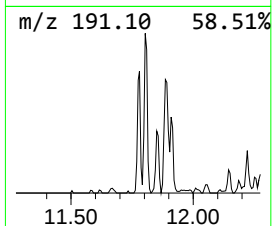
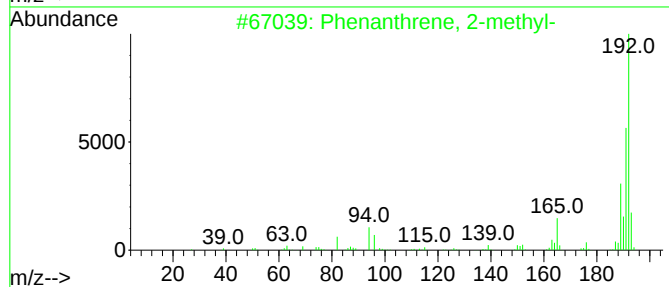
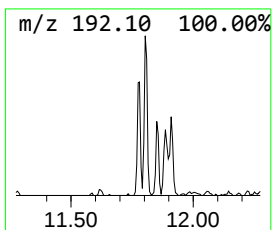
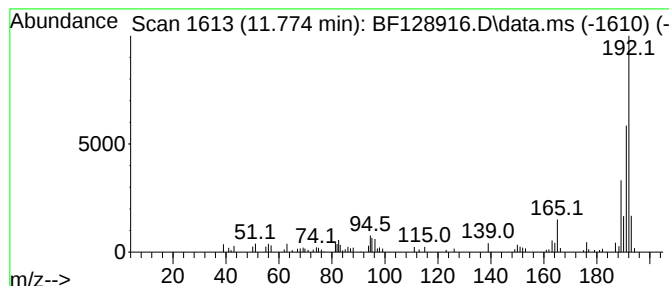
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF060222.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.774	2.05 ng	64602	Phenanthrene-d10	11.275

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	97
3		1H-Cyclopropa[l]phenanthrene,1a,...	192	C15H12	000949-41-7	95
4		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	95
5		Anthracene, 9-methyl-	192	C15H12	000779-02-2	93



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Sample : N3421-04
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
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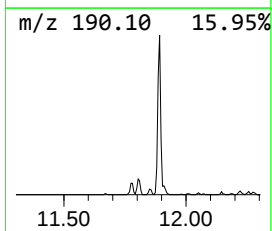
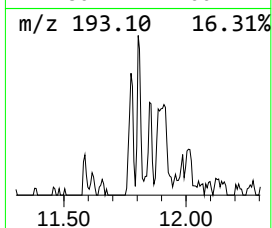
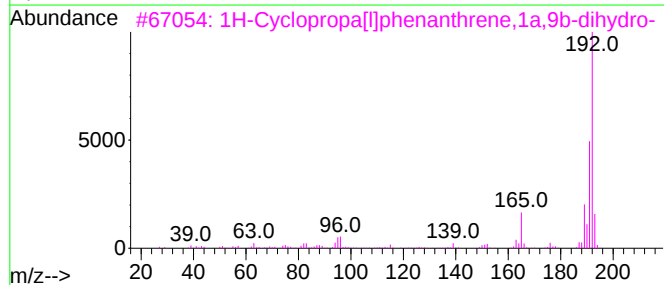
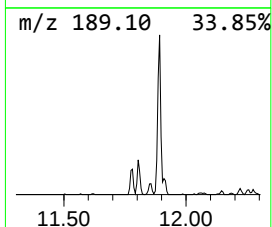
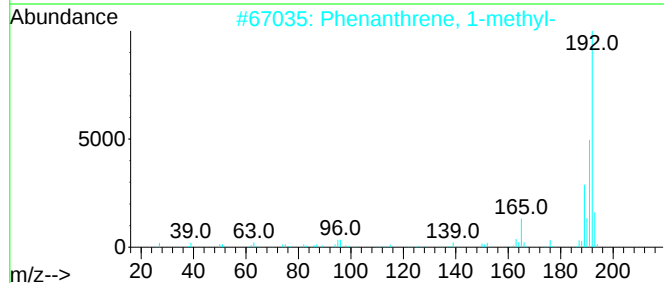
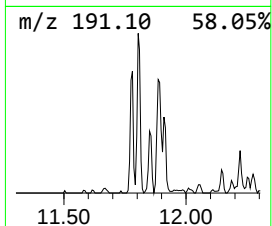
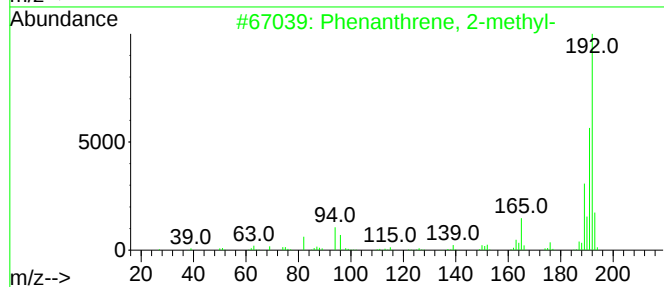
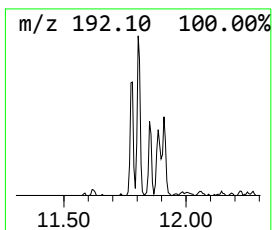
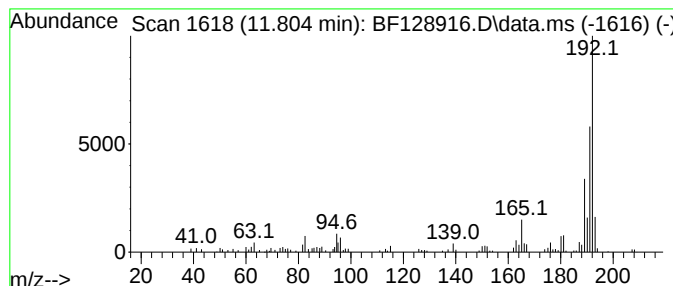
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF060222.M
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.804	3.03 ng	95320	Phenanthrene-d10	11.275

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	97
3		1H-Cyclopropa[1]phenanthrene,1a,...	192	C15H12	000949-41-7	96
4		Anthracene, 1-methyl-	192	C15H12	000610-48-0	94
5		Anthracene, 2-methyl-	192	C15H12	000613-12-7	93



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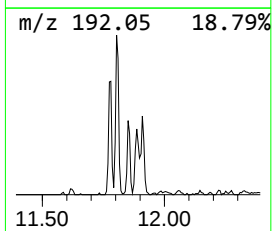
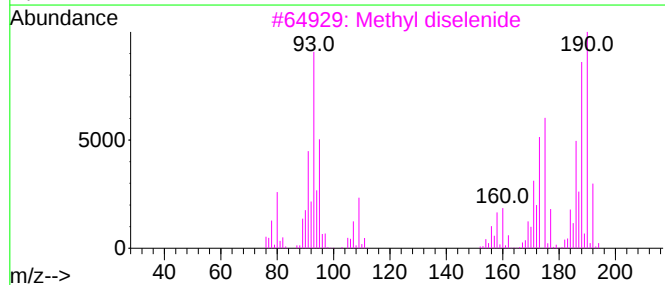
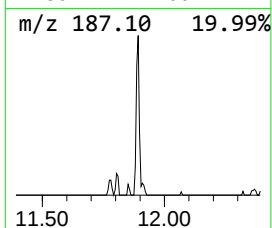
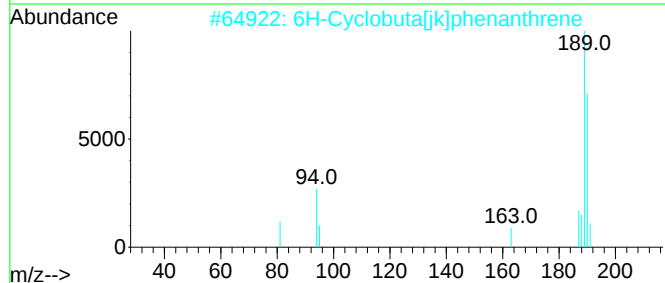
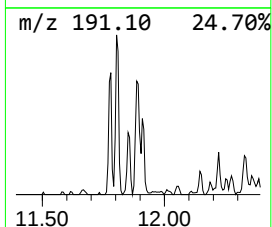
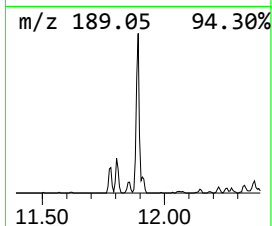
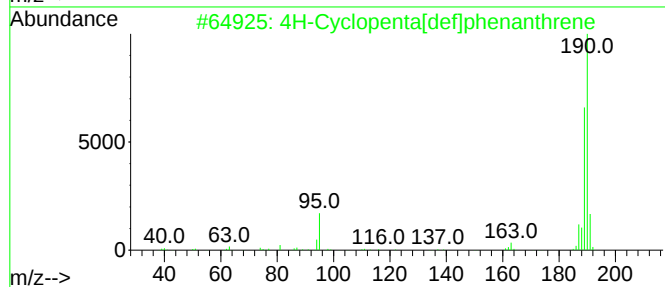
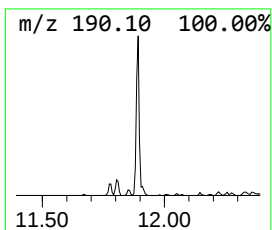
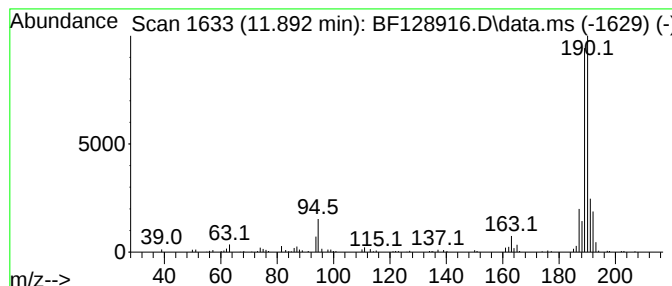
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF060222.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 4H-Cyclopenta[def]phenanthrene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD		R.T.	
11.892	4.89 ng	153763	Phenanthrene-d10		11.275	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene		190	C15H10	000203-64-5	87
2	6H-Cyclobuta[jk]phenanthrene		190	C15H10	083469-43-6	56
3	Methyl diselenide		190	C2H6Se2	007101-31-7	49
4	2,2'-Bis(4,5-dimethylimidazole)		190	C10H14N4	069286-06-2	45
5	1,4-Naphthalenedione, 2,5-dihydr...		190	C10H6O4	004923-55-1	32



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SB-3-0-2-20220620

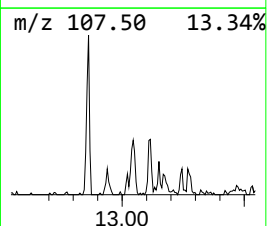
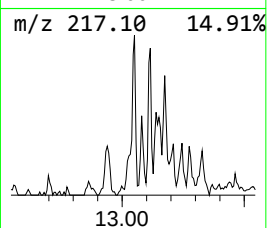
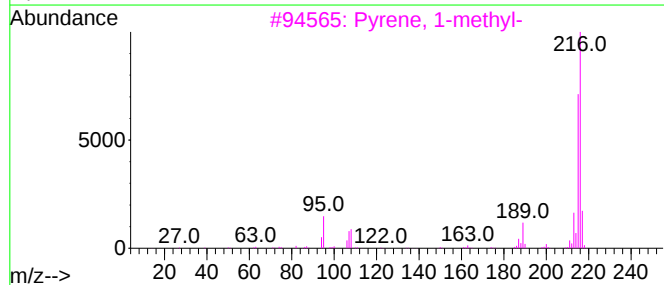
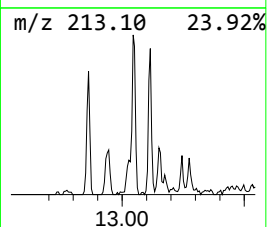
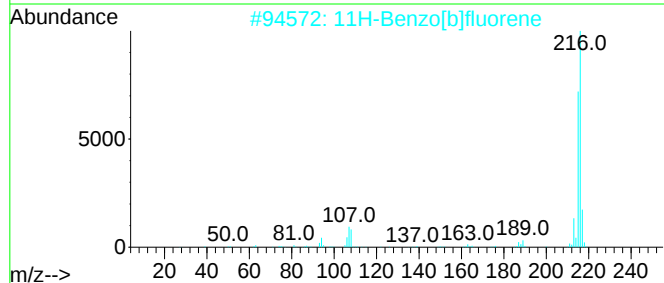
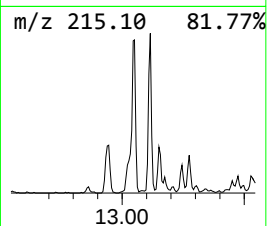
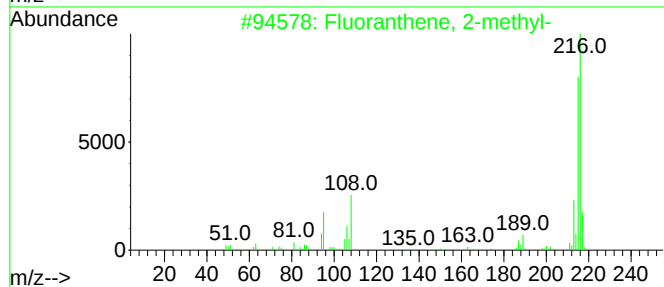
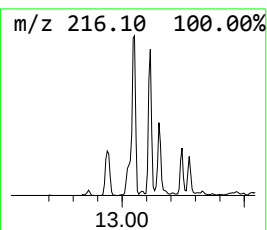
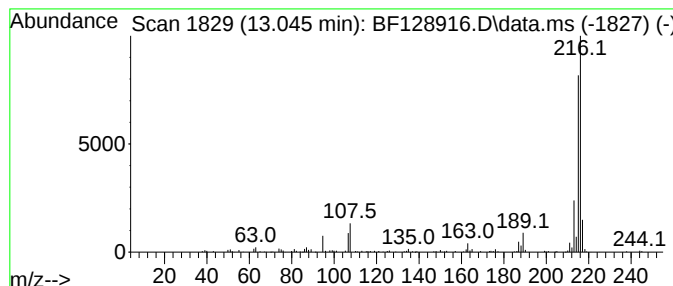
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 4 Fluoranthene, 2-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.045	2.38 ng	165748	Chrysene-d12	13.921

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	94
2			11H-Benzo[b]fluorene	216	C17H12	000243-17-4	91
3			Pyrene, 1-methyl-	216	C17H12	002381-21-7	90
4			11H-Benzo[a]fluorene	216	C17H12	000238-84-6	87
5			7H-Benzo[c]fluorene	216	C17H12	000205-12-9	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF062122\
Data File : BF128916.D
Acq On : 22 Jun 2022 00:19
Operator : CG\JU
Sample : N3421-04
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-3-0-2-20220620

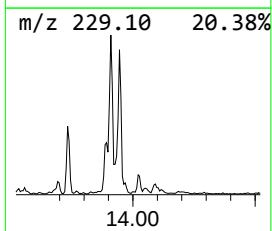
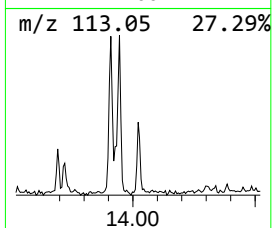
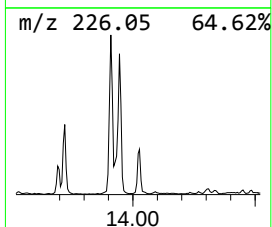
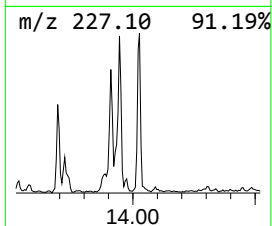
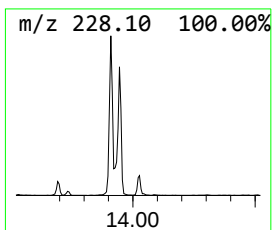
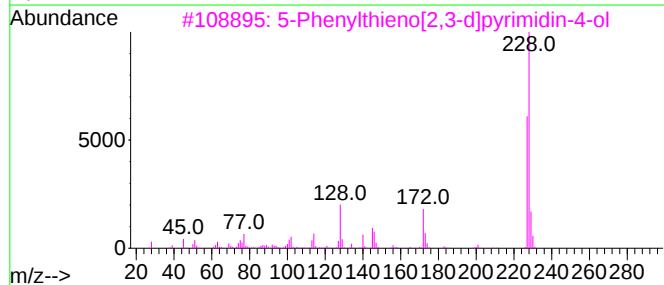
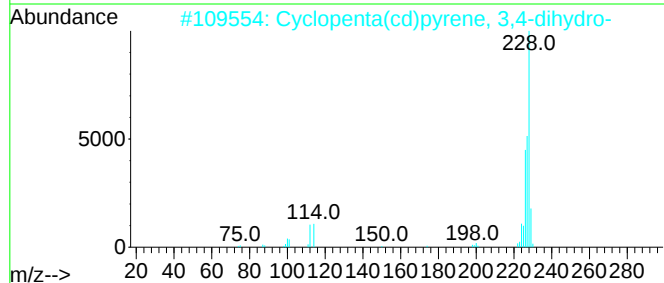
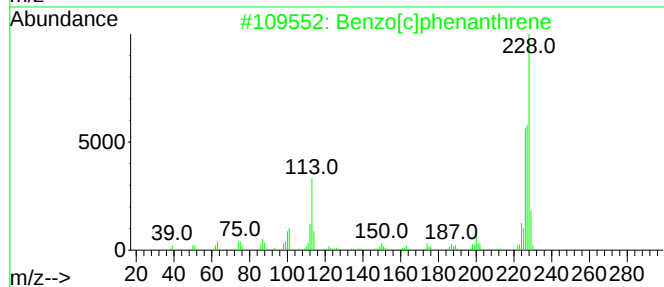
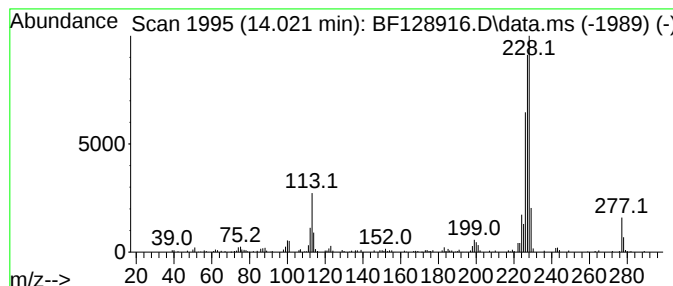
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Benzo[c]phenanthrene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.021	2.84 ng	198164	Chrysene-d12	13.921

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[c]phenanthrene	228	C18H12	000195-19-7	60
2			Cyclopenta(cd)pyrene, 3,4-dihydro-	228	C18H12	025732-74-5	55
3			5-Phenylthieno[2,3-d]pyrimidin-4-ol	228	C12H8N2O5	035978-39-3	52
4			Triphenylene	228	C18H12	000217-59-4	46
5			Isothiazol-5-amine, 4-bromo-3-ch...	226	C4H4BrClN2S	1000272-59-9	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF062122\
Data File : BF128916.D
Acq On : 22 Jun 2022 00:19
Operator : CG\JU
Sample : N3421-04
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
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ClientSampleId :
SB-3-0-2-20220620

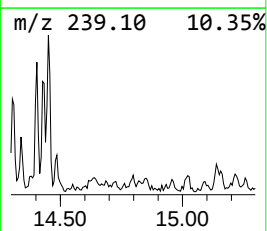
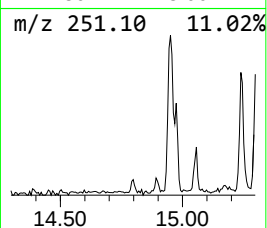
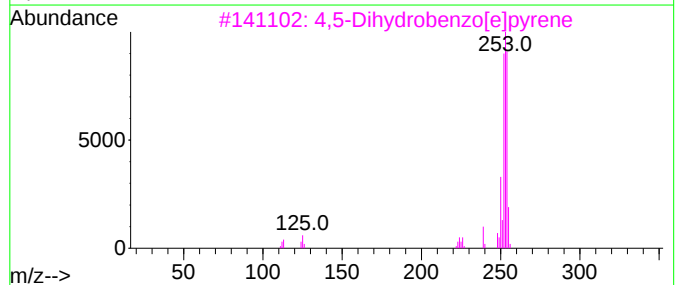
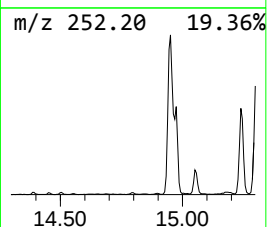
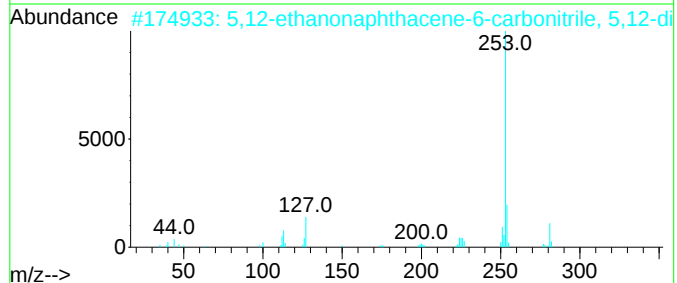
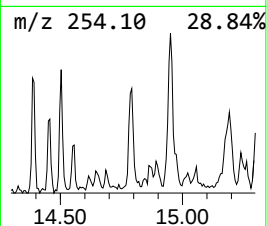
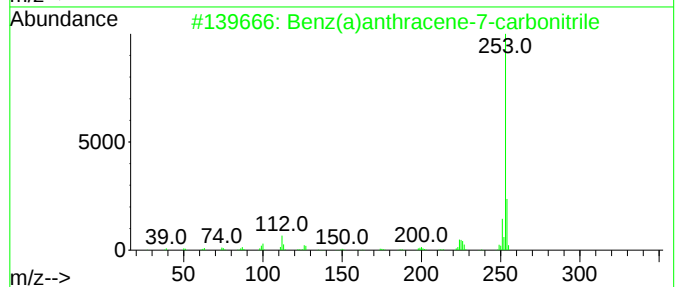
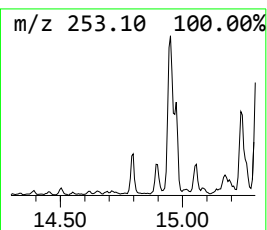
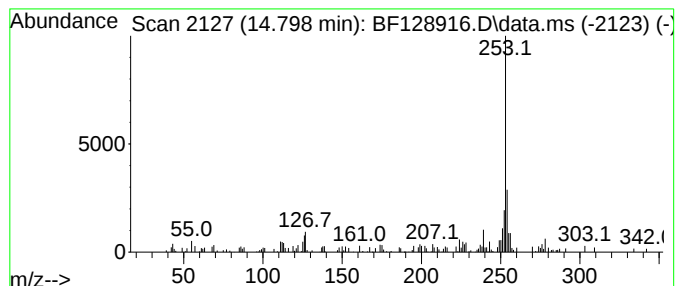
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Benz(a)anthracene-7-carboni... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.798	3.24 ng	65506	Perylene-d12	15.362

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benz(a)anthracene-7-carbonitrile	253	C19H11N	007476-08-6	53
2			5,12-ethanonaphthacene-6-carboni...	281	C21H15N	1000397-19-0	53
3			4,5-Dihydrobenzo[e]pyrene	254	C20H14	095676-42-9	50
4			2-Methyl-4-trihexylsilyloxyoct-5...	422	C27H54OSi	1000299-52-5	47
5			Fumaric acid, 3-hexyl undecyl ester	354	C21H38O4	1000348-61-0	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF062122\
Data File : BF128916.D
Acq On : 22 Jun 2022 00:19
Operator : CG\JU
Sample : N3421-04
Misc :
ALS Vial : 17 Sample Multiplier: 1

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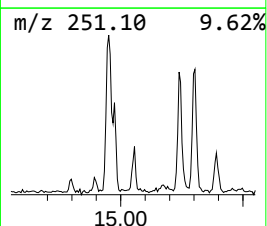
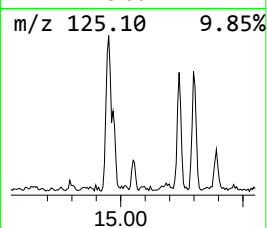
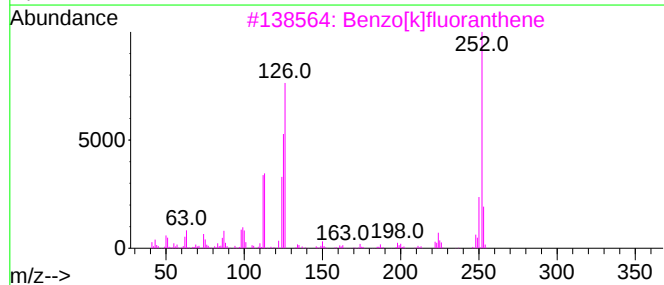
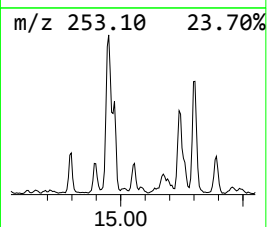
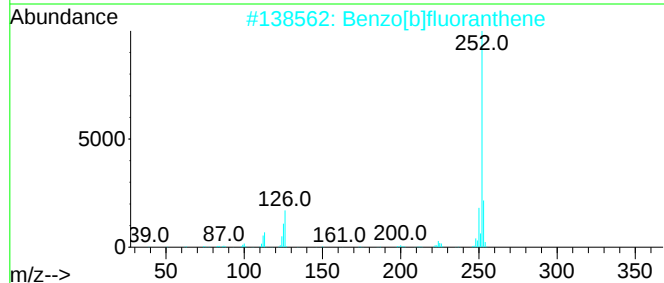
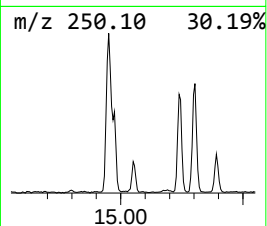
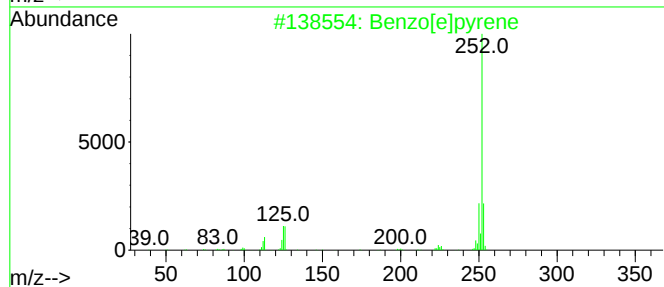
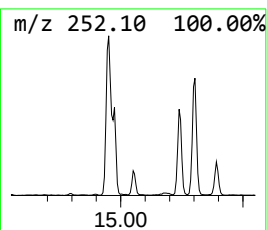
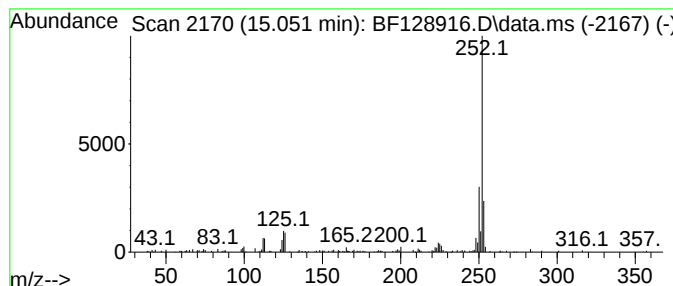
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Benzo[e]pyrene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.051	3.39 ng	68639	Perylene-d12	15.362

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[e]pyrene	252	C20H12	000192-97-2	97
2			Benzo[b]fluoranthene	252	C20H12	000205-99-2	93
3			Benzo[k]fluoranthene	252	C20H12	000207-08-9	93
4			Benzo[a]pyrene	252	C20H12	000050-32-8	91
5			Perylene	252	C20H12	000198-55-0	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF062122\
Data File : BF128916.D
Acq On : 22 Jun 2022 00:19
Operator : CG\JU
Sample : N3421-04
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF060222.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...	11.774	2.0	ng	64602	4	11.275	629211	20.0
Phenanthrene, 1...	11.804	3.0	ng	95320	4	11.275	629211	20.0
4H-Cyclopenta[d...	11.892	4.9	ng	153763	4	11.275	629211	20.0
Fluoranthene, 2...	13.045	2.4	ng	165748	5	13.921	1394430	20.0
Benzo[c]phenant...	14.021	2.8	ng	198164	5	13.921	1394430	20.0
Benz(a)anthrace...	14.798	3.2	ng	65506	6	15.362	404355	20.0
Benzo[e]pyrene	15.051	3.4	ng	68639	6	15.362	404355	20.0