

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF101520\
 Data File : BF122002.D
 Acq On : 15 Oct 2020 20:51
 Operator : JU/CG
 Sample : L4406-01 2X
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 JC-02-101420-A

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-BF101220.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF122002.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.357	553	556	575	rBV	847651	1056636	78.63%	5.280%
2	6.387	727	731	757	rBV	904448	1136117	84.55%	5.678%
3	6.739	788	791	798	rBV	719119	632489	47.07%	3.161%
4	7.310	884	888	904	rBV	703499	760762	56.61%	3.802%
5	8.028	1006	1010	1023	rBV	822848	845099	62.89%	4.223%
6	9.098	1188	1192	1202	rBV	1541679	1343748	100.00%	6.715%
7	9.498	1254	1260	1268	rVB	155532	159812	11.89%	0.799%
8	9.639	1281	1284	1288	rVB	17360	16881	1.26%	0.084%
9	9.775	1303	1307	1311	rBV	1014224	918436	68.35%	4.590%
10	9.810	1311	1313	1321	rVB	87040	90515	6.74%	0.452%
11	9.986	1340	1343	1346	rVB	31372	28024	2.09%	0.140%
12	10.327	1397	1401	1406	rBV	59744	63116	4.70%	0.315%
13	10.480	1425	1427	1430	rVB	21372	17684	1.32%	0.088%
14	10.569	1438	1442	1453	rBV	572023	617629	45.96%	3.087%
15	10.680	1458	1461	1469	rVB4	23242	43371	3.23%	0.217%
16	11.075	1524	1528	1532	rVV	46986	44071	3.28%	0.220%
17	11.122	1532	1536	1539	rVV3	43210	54964	4.09%	0.275%
18	11.157	1539	1542	1549	rVB	62453	64835	4.82%	0.324%
19	11.263	1556	1560	1562	rBV	904364	875179	65.13%	4.374%
20	11.286	1562	1564	1568	rVV	898116	734599	54.67%	3.671%
21	11.339	1569	1573	1576	rVV	129509	130327	9.70%	0.651%
22	11.380	1578	1580	1585	rVB2	14311	18986	1.41%	0.095%
23	11.498	1596	1600	1606	rBV2	110087	134147	9.98%	0.670%
24	11.551	1606	1609	1611	rVV2	33069	29003	2.16%	0.145%
25	11.604	1614	1618	1621	rVB3	15132	20026	1.49%	0.100%
26	11.651	1623	1626	1629	rBV	89826	70503	5.25%	0.352%
27	11.763	1642	1645	1648	rBV	94566	81826	6.09%	0.409%
28	11.792	1648	1650	1654	rVB	129741	104158	7.75%	0.521%
29	11.839	1655	1658	1661	rVV	25677	21547	1.60%	0.108%
30	11.874	1661	1664	1666	rVV	136793	136613	10.17%	0.683%
31	11.898	1666	1668	1671	rVB	69614	55515	4.13%	0.277%
32	11.957	1674	1678	1680	rBV3	23610	28926	2.15%	0.145%
33	12.057	1692	1695	1698	rBV	69962	58196	4.33%	0.291%
34	12.092	1698	1701	1706	rVV	90938	86026	6.40%	0.430%
35	12.204	1718	1720	1724	rVB3	20941	15737	1.17%	0.079%
36	12.239	1724	1726	1728	rBV2	28813	21072	1.57%	0.105%

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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-BF101220.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

37	12.310	1735	1738	1740	rBV	52230	53519	3.98%	0.267%
38	12.333	1740	1742	1744	rVV	190511	185877	13.83%	0.929%
39	12.357	1744	1746	1751	rVV2	239234	248230	18.47%	1.241%
40	12.404	1751	1754	1758	rVV2	72364	88810	6.61%	0.444%

41	12.445	1758	1761	1763	rVV2	41052	48598	3.62%	0.243%
42	12.474	1763	1766	1770	rVB	1181854	1000878	74.48%	5.002%
43	12.539	1775	1777	1780	rVB	25905	27297	2.03%	0.136%
44	12.621	1788	1791	1794	rVB3	48282	50858	3.78%	0.254%
45	12.704	1799	1805	1811	rBV	945708	982522	73.12%	4.910%

46	12.751	1811	1813	1819	rVB2	44920	54042	4.02%	0.270%
47	12.845	1822	1829	1832	rBV	1204234	1119150	83.29%	5.593%
48	12.869	1832	1833	1838	rVB	49970	42287	3.15%	0.211%
49	12.927	1838	1843	1846	rBV3	64247	111277	8.28%	0.556%
50	12.957	1846	1848	1850	rBV	23276	19267	1.43%	0.096%

51	13.033	1853	1861	1864	rBV2	115108	203808	15.17%	1.019%
52	13.074	1865	1868	1869	rBV3	23769	25773	1.92%	0.129%
53	13.098	1869	1872	1874	rVB2	48898	40058	2.98%	0.200%
54	13.133	1874	1878	1881	rBV2	83318	102215	7.61%	0.511%
55	13.227	1891	1894	1897	rBV	48052	47453	3.53%	0.237%

56	13.263	1897	1900	1902	rBV2	48796	60533	4.50%	0.303%
57	13.416	1923	1926	1928	rBV4	23834	27300	2.03%	0.136%
58	13.457	1931	1933	1935	rVV	25427	21157	1.57%	0.106%
59	13.545	1945	1948	1952	rVB	112454	97486	7.25%	0.487%
60	13.651	1963	1966	1968	rBV	106587	96991	7.22%	0.485%

61	13.674	1968	1970	1973	rVB	69671	61266	4.56%	0.306%
62	13.704	1973	1975	1977	rBV	65872	68865	5.12%	0.344%
63	13.757	1981	1984	1987	rVV	110246	95565	7.11%	0.478%
64	13.904	2001	2009	2011	rBV2	934927	1277080	95.04%	6.382%
65	13.927	2011	2013	2016	rVB	453366	369556	27.50%	1.847%

66	14.010	2024	2027	2029	rVB	61439	56641	4.22%	0.283%
67	14.063	2033	2036	2040	rBV4	34822	55506	4.13%	0.277%
68	14.139	2047	2049	2051	rVB	41714	33161	2.47%	0.166%
69	14.292	2072	2075	2077	rVB	87935	81746	6.08%	0.409%
70	14.321	2078	2080	2082	rBV	42370	32924	2.45%	0.165%

71	14.774	2155	2157	2160	rBV2	54389	58407	4.35%	0.292%
72	14.933	2179	2184	2186	rBV	395606	566449	42.15%	2.831%
73	15.033	2198	2201	2212	rVB	79340	117066	8.71%	0.585%
74	15.221	2229	2233	2238	rVB	237850	287306	21.38%	1.436%
75	15.280	2239	2243	2248	rVB	328559	382578	28.47%	1.912%

76	15.339	2249	2253	2257	rBV	566448	719816	53.57%	3.597%
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ClientSampleId :
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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	16.751	2488	2493	2502	rVB	132376	271941	20.24%	1.359%
78	17.180	2561	2566	2575	rVB	100203	202350	15.06%	1.011%

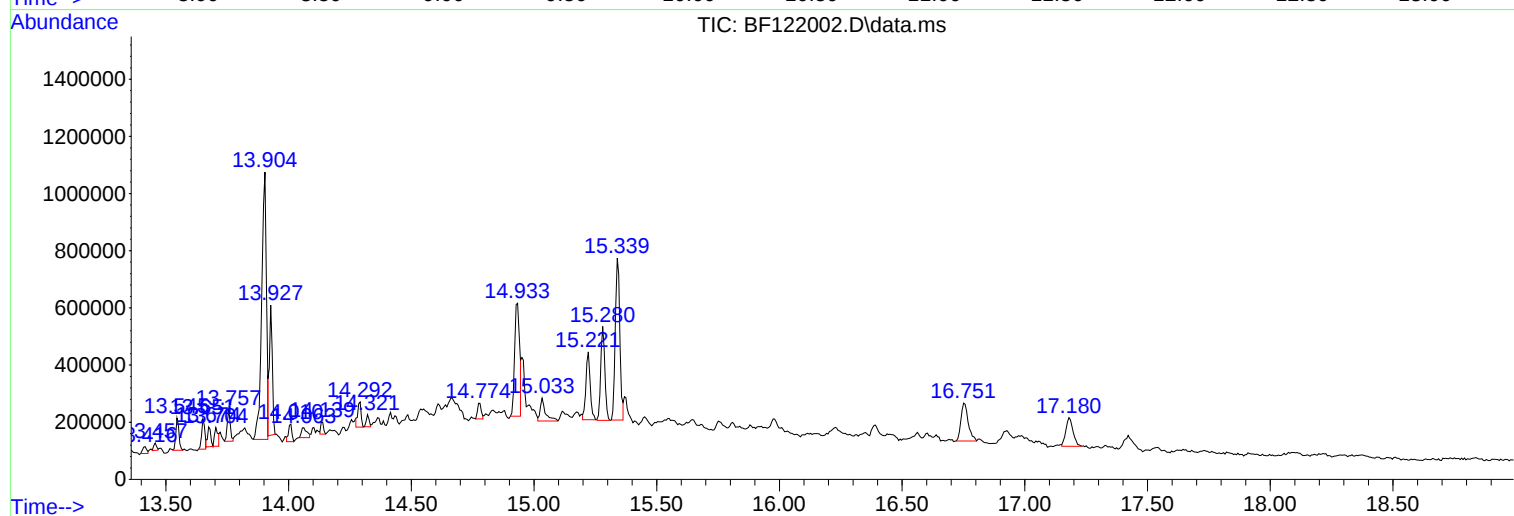
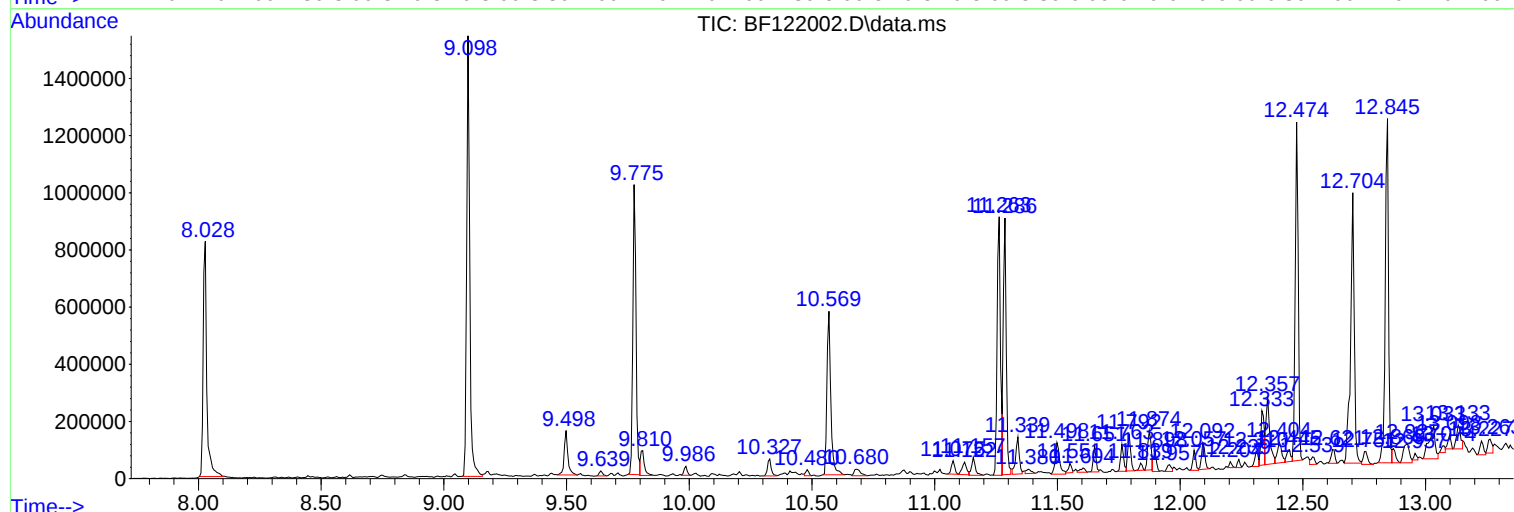
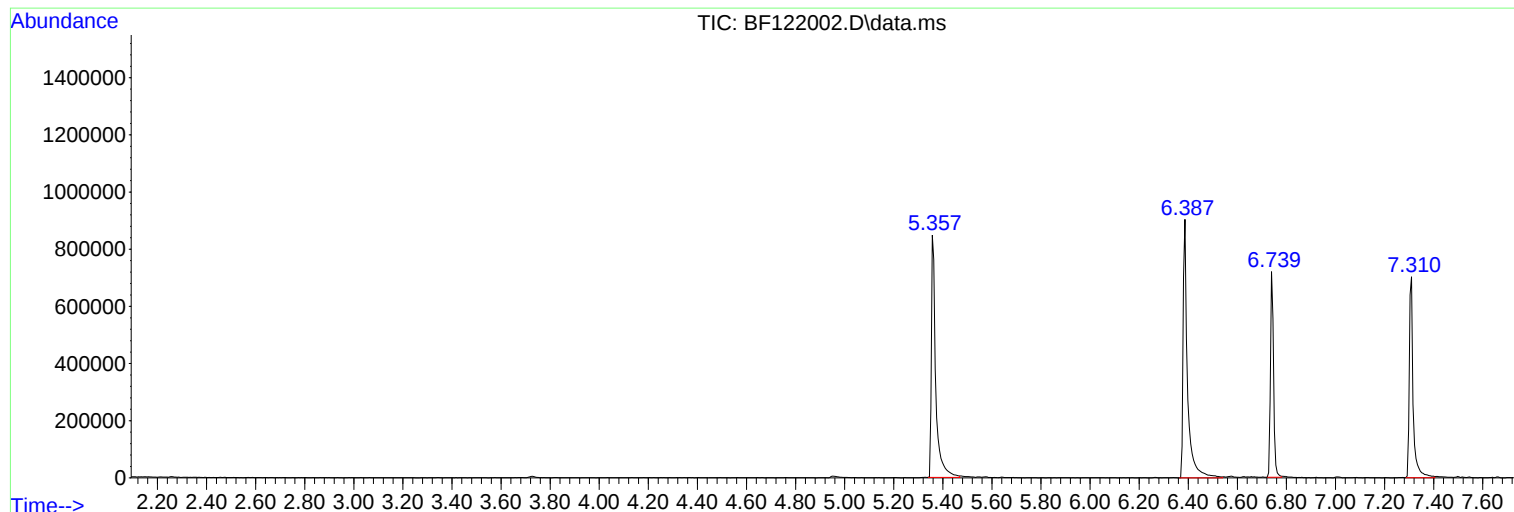
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Misc :
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Instrument :
BNA_F
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF101220.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF101520\
Data File : BF122002.D
Acq On : 15 Oct 2020 20:51
Operator : JU/CG
Sample : L4406-01 2X
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
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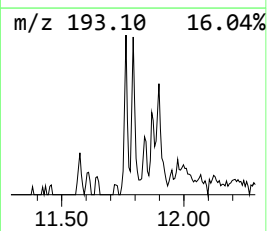
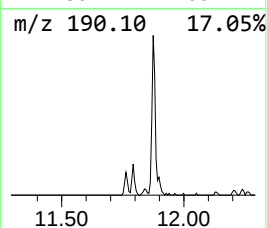
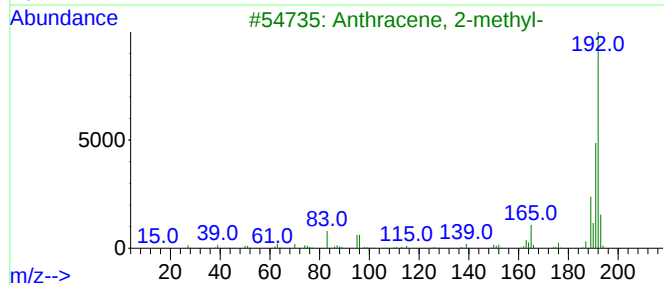
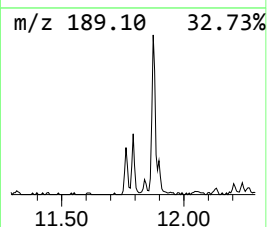
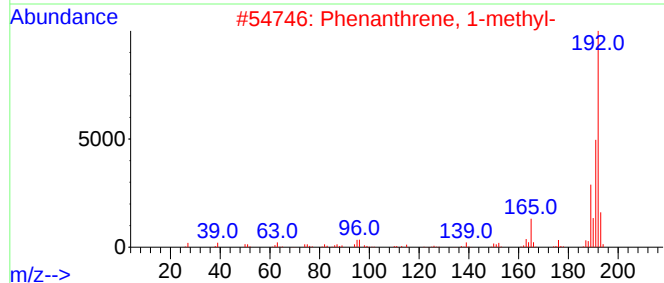
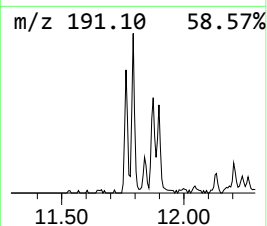
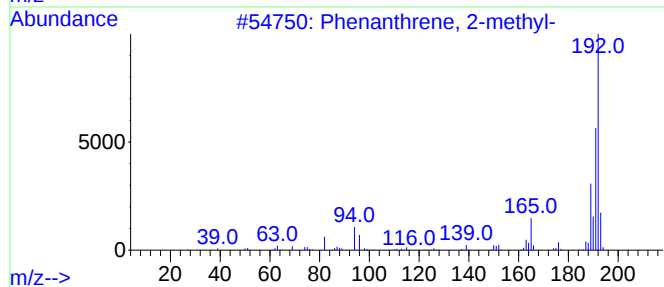
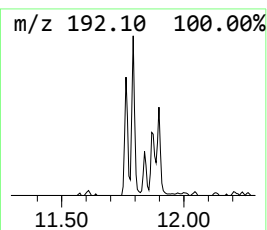
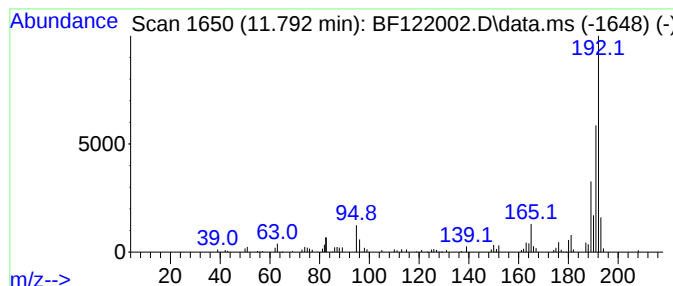
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF101220.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Phenanthrene, 2-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.792	2.38 ng	104158	Phenanthrene-d10	11.263

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	97
2			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	97
3			Anthracene, 2-methyl-	192	C15H12	000613-12-7	96
4			Anthracene, 1-methyl-	192	C15H12	000610-48-0	93
5			1H-Indene, 1-phenyl-	192	C15H12	001961-96-2	93



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Misc :
ALS Vial : 20 Sample Multiplier: 1

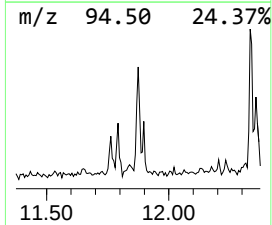
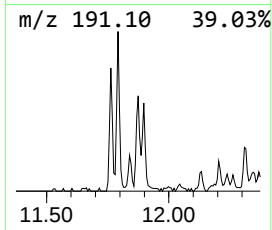
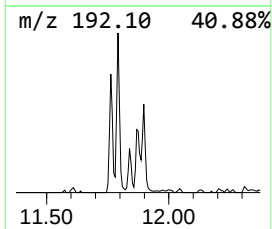
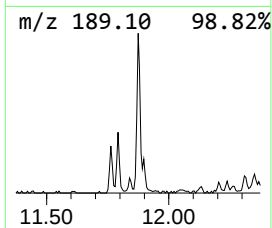
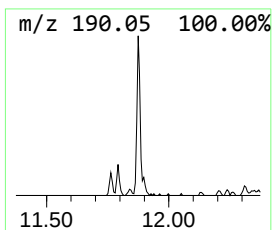
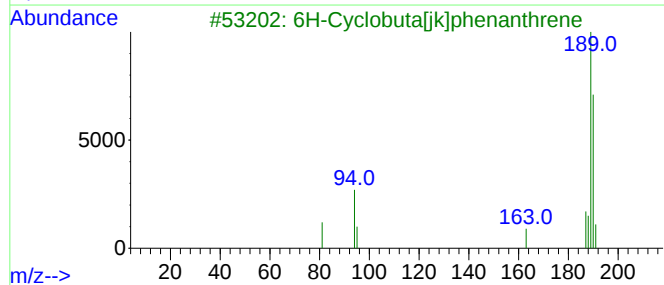
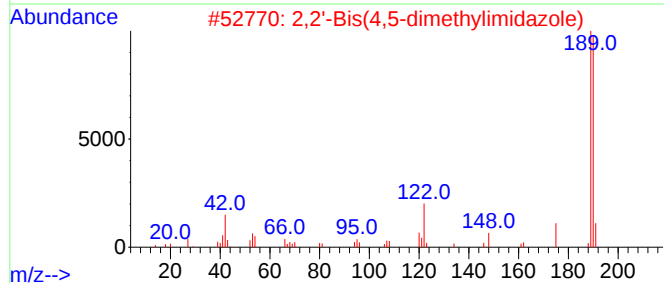
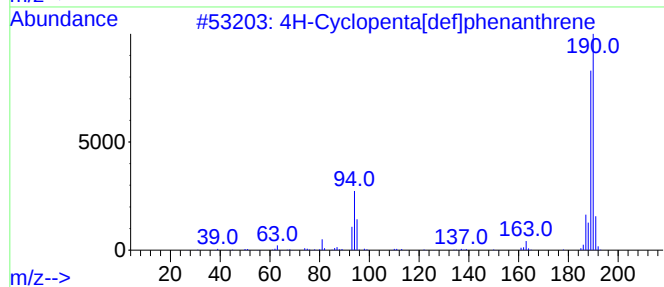
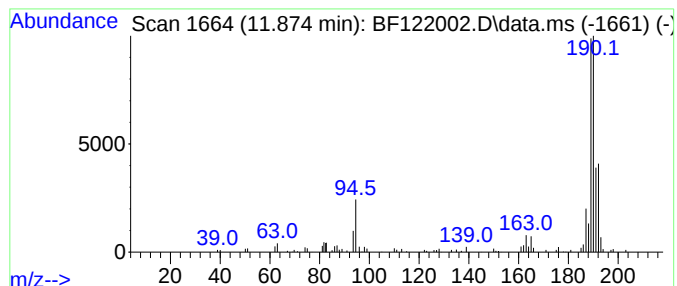
Instrument :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF101220.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 4H-Cyclopenta[def]phenanthrene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD		R.T.	
11.874	3.12 ng	136613	Phenanthrene-d10		11.263	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene		190	C15H10	000203-64-5	76
2	2,2'-Bis(4,5-dimethylimidazole)		190	C10H14N4	069286-06-2	37
3	6H-Cyclobuta[jk]phenanthrene		190	C15H10	083469-43-6	28
4	1,1'-Biphenyl, 4,4'-difluoro-		190	C12H8F2	000398-23-2	23
5	Cyclohexanone, 2-(2-furanylmethy...		190	C12H14O2	056053-07-7	14



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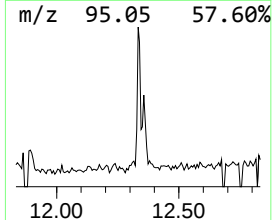
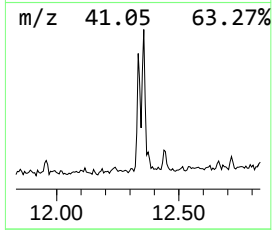
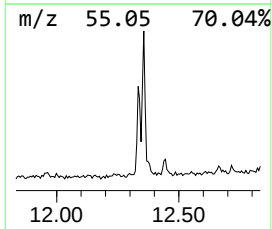
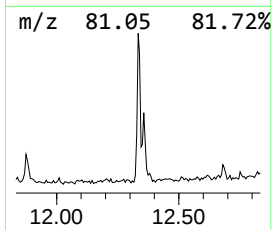
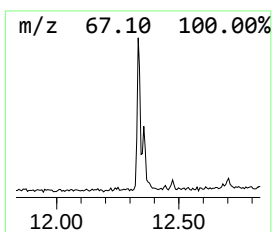
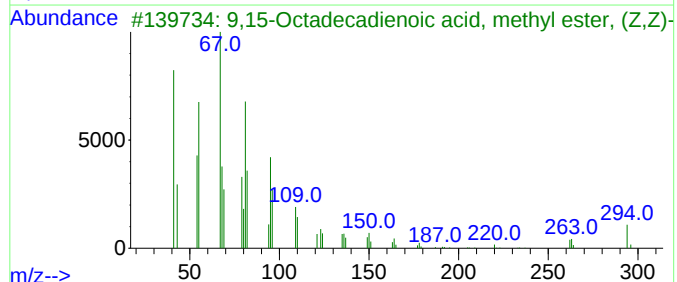
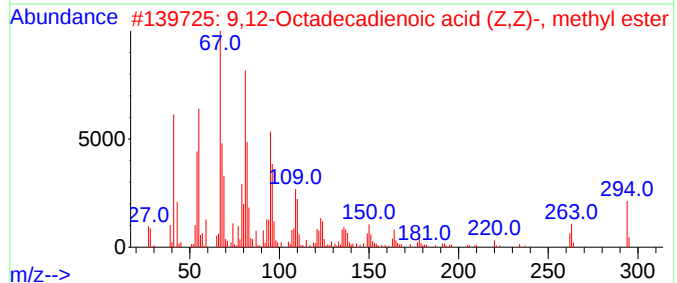
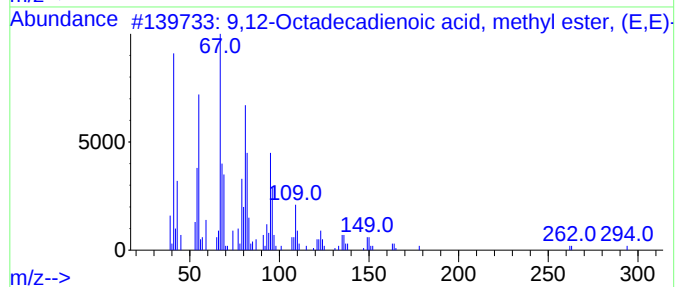
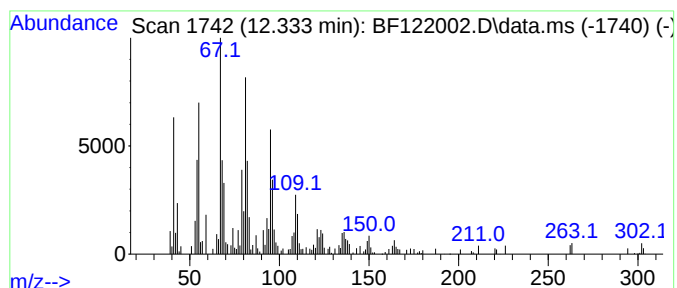
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF101220.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 9,12-Octadecadienoic acid, ... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.333	4.25 ng	185877	Phenanthrene-d10	11.263

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	9,12-Octadecadienoic acid, methy...	294	C19H34O2	002566-97-4	99		
2	9,12-Octadecadienoic acid (Z,Z)-...	294	C19H34O2	000112-63-0	99		
3	9,15-Octadecadienoic acid, methy...	294	C19H34O2	017309-05-6	99		
4	8,11-Octadecadienoic acid, methy...	294	C19H34O2	056599-58-7	99		
5	10,13-Octadecadienoic acid, meth...	294	C19H34O2	056554-62-2	99		



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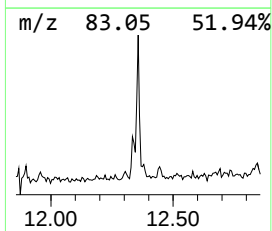
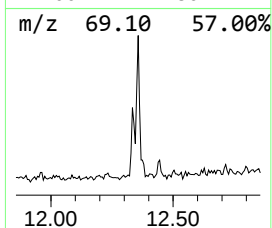
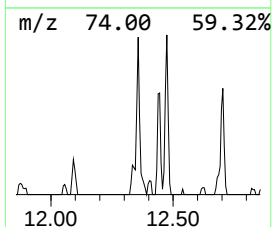
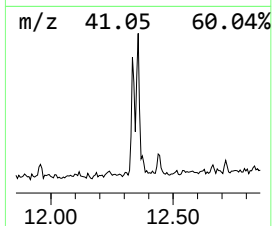
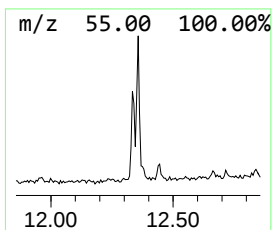
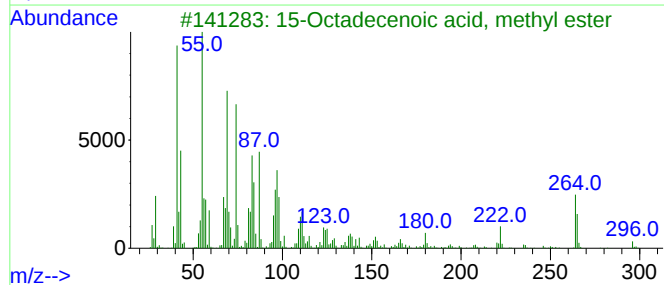
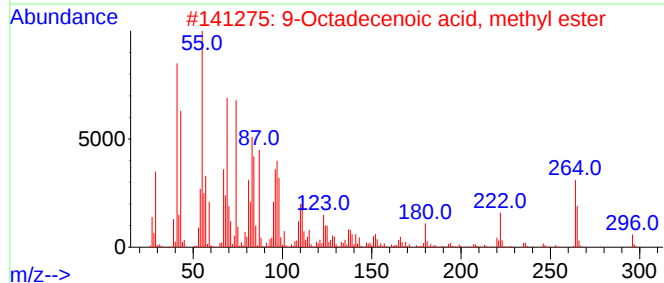
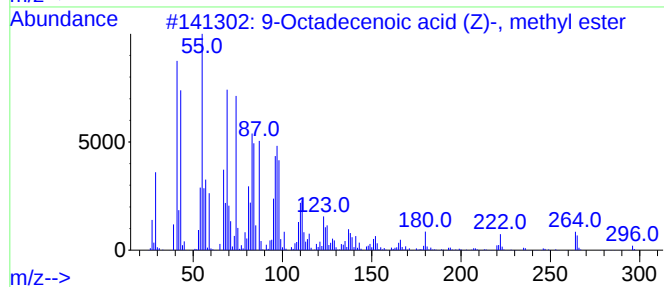
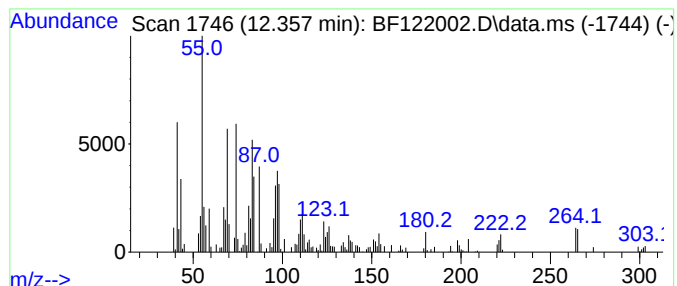
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 9-Octadecenoic acid (Z)-, m... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.357	5.67 ng	248230	Phenanthrene-d10	11.263

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			9-Octadecenoic acid (Z)-, methyl...	296	C19H36O2	000112-62-9	95
2			9-Octadecenoic acid, methyl ester	296	C19H36O2	002462-84-2	94
3			15-Octadecenoic acid, methyl ester	296	C19H36O2	004764-72-1	91
4			13-Octadecenoic acid, methyl ester	296	C19H36O2	056554-47-3	91
5			11-Octadecenoic acid, methyl ester	296	C19H36O2	052380-33-3	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF101520\
Data File : BF122002.D
Acq On : 15 Oct 2020 20:51
Operator : JU/CG
Sample : L4406-01 2X
Misc :
ALS Vial : 20 Sample Multiplier: 1

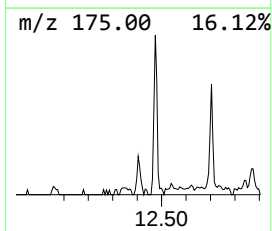
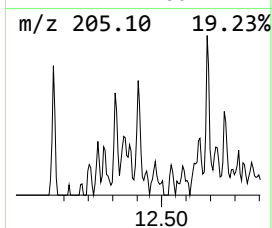
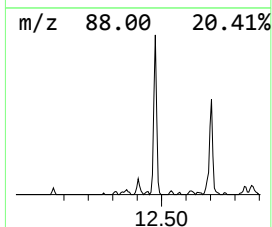
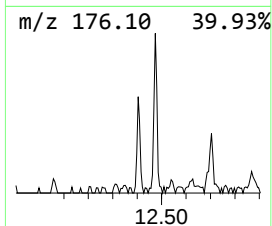
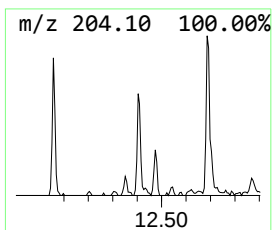
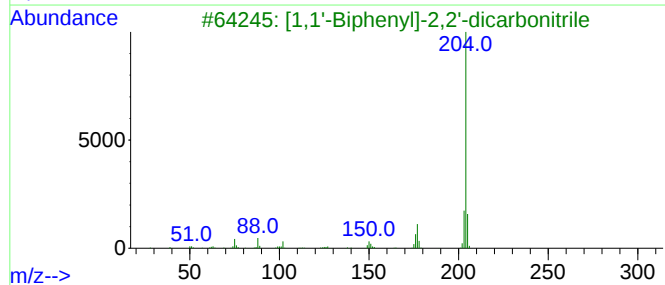
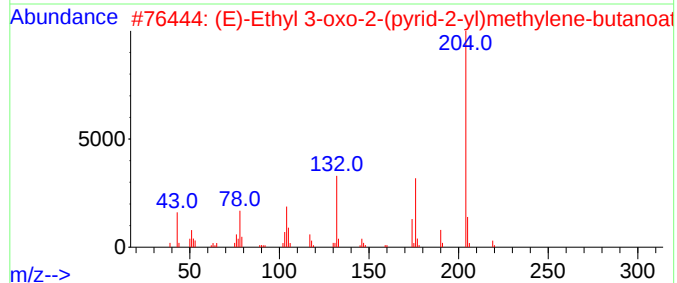
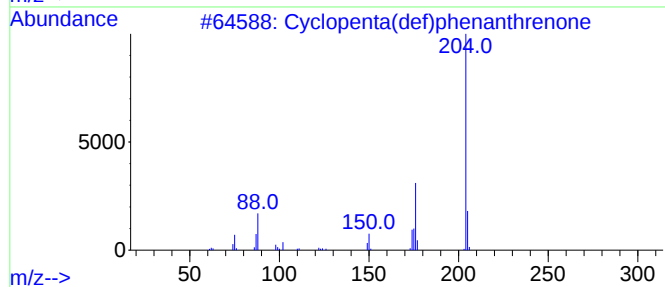
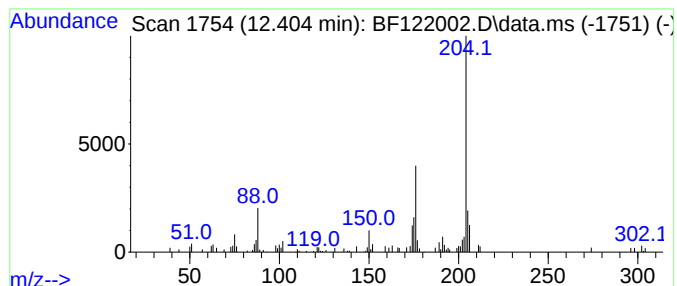
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ClientSampleId :
JC-02-101420-A

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF101220.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Cyclopenta(def)phenanthrenone Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.404	2.03 ng	88810	Phenanthrene-d10	11.263	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopenta(def)phenanthrenone	204	C15H8O	005737-13-3	93
2	(E)-Ethyl 3-oxo-2-(pyrid-2-yl)me...	219	C12H13NO3	1000147-59-7	45
3	[1,1'-Biphenyl]-2,2'-dicarbonitrile	204	C14H8N2	004341-02-0	43
4	4(equat)-(but-3-en-1-ynyl)-2(axi...	219	C14H21NO	038423-22-2	42
5	N-(4-Chloro-phenyl)-2-(3,4-dihyd...	330	C17H15ClN2OS	1000296-34-3	42



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF101520\
Data File : BF122002.D
Acq On : 15 Oct 2020 20:51
Operator : JU/CG
Sample : L4406-01 2X
Misc :
ALS Vial : 20 Sample Multiplier: 1

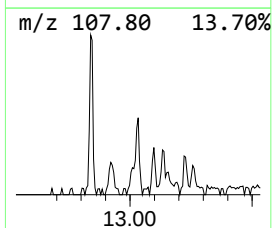
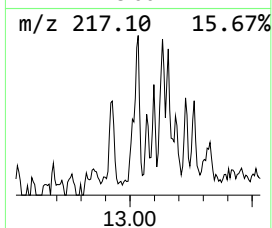
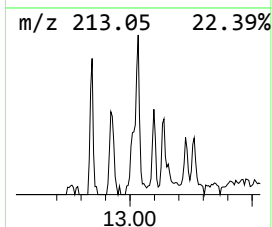
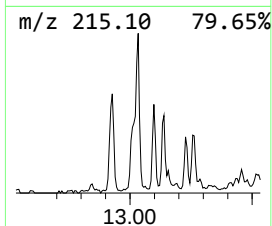
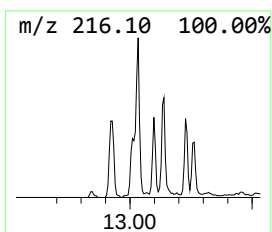
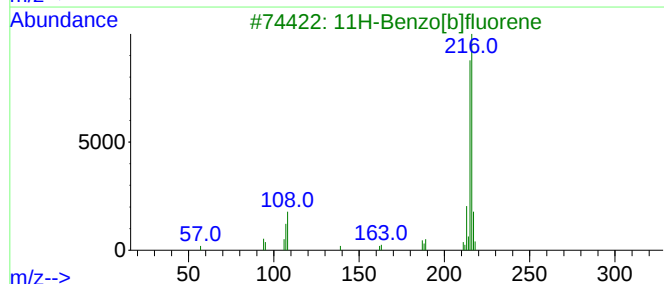
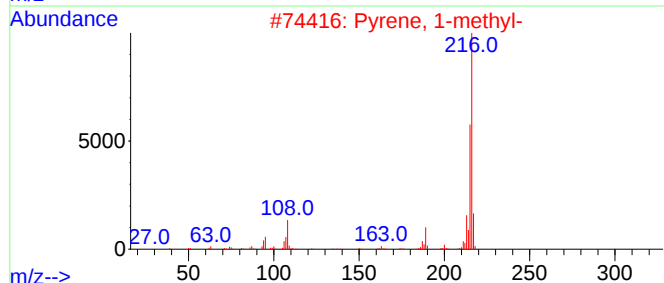
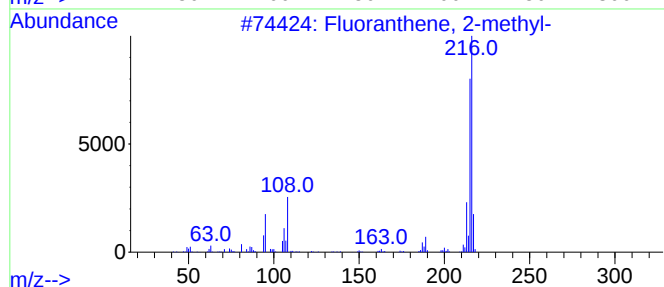
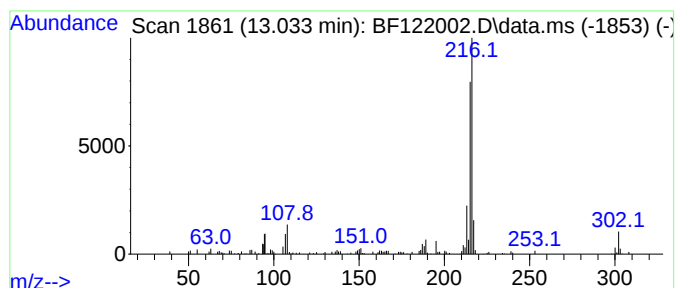
Instrument :
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ClientSampleId :
JC-02-101420-A

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF101220.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Fluoranthene, 2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.033	3.19 ng	203808	Chrysene-d12	13.904	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	93
2	Pyrene, 1-methyl-	216	C17H12	002381-21-7	93
3	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	91
4	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	90
5	Pyrene, 2-methyl-	216	C17H12	003442-78-2	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF101520\
Data File : BF122002.D
Acq On : 15 Oct 2020 20:51
Operator : JU/CG
Sample : L4406-01 2X
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
JC-02-101420-A

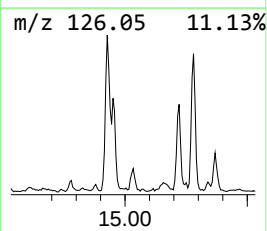
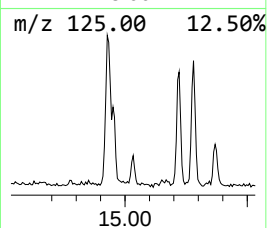
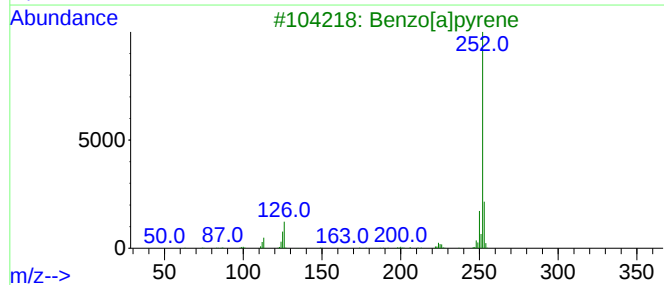
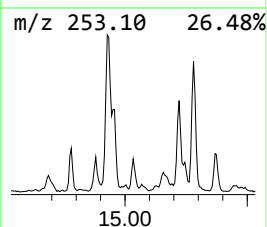
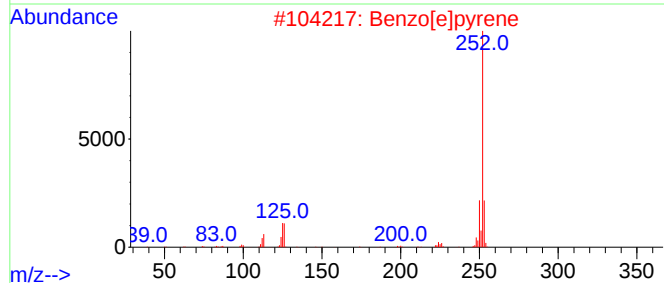
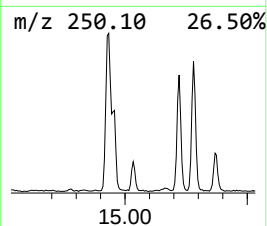
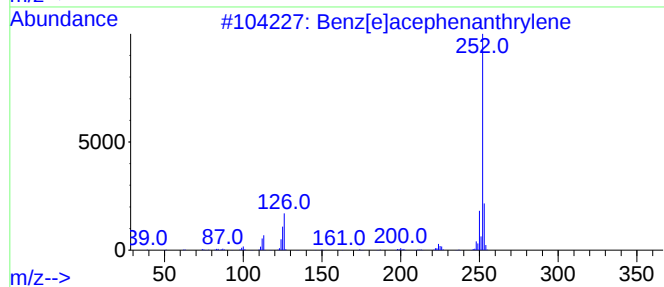
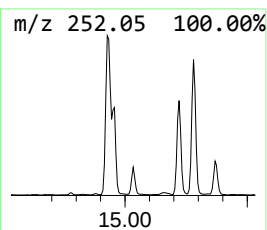
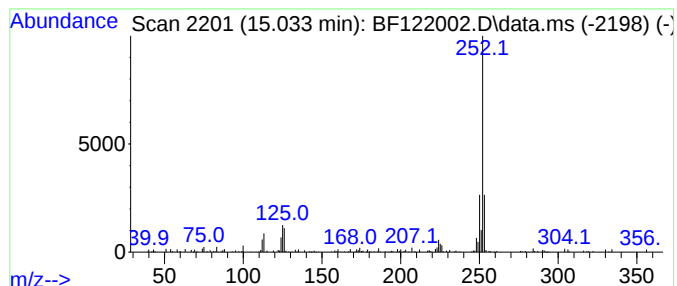
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Benzo[e]pyrene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.033	3.25 ng	117066	Perylene-d12	15.339

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benz[e]acephenanthrylene	252	C20H12	000205-99-2	93
2			Benzo[e]pyrene	252	C20H12	000192-97-2	91
3			Benzo[a]pyrene	252	C20H12	000050-32-8	90
4			Benzo[k]fluoranthene	252	C20H12	000207-08-9	89
5			Perylene	252	C20H12	000198-55-0	81



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF101520\
Data File : BF122002.D
Acq On : 15 Oct 2020 20:51
Operator : JU/CG
Sample : L4406-01 2X
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
JC-02-101420-A

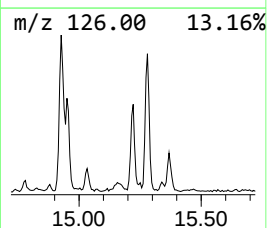
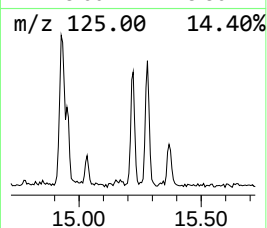
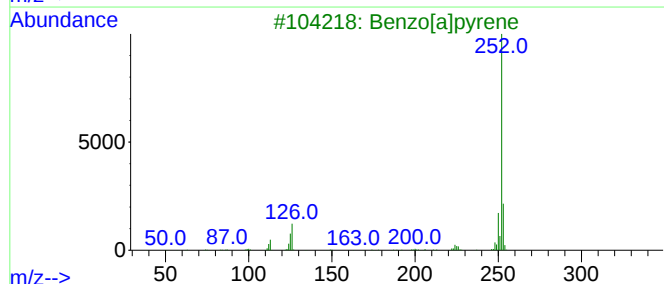
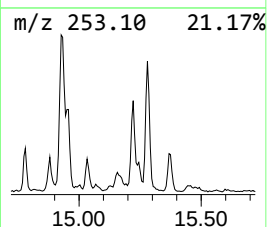
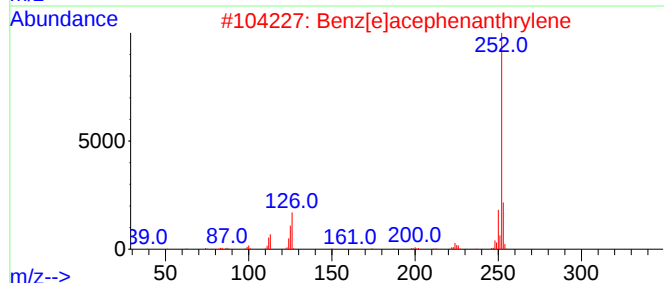
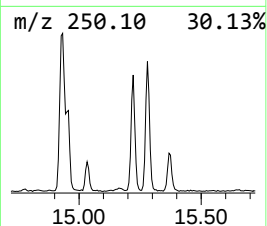
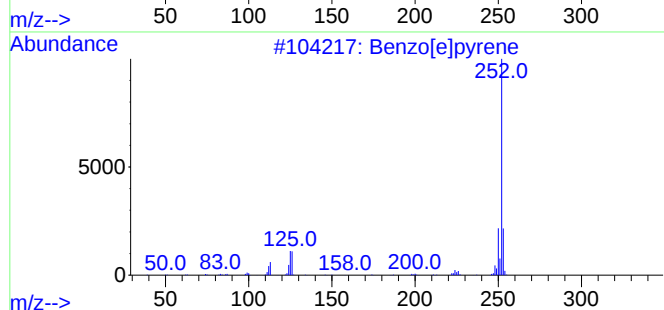
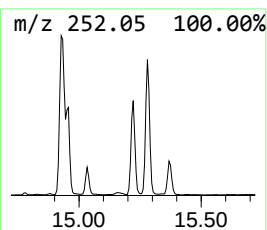
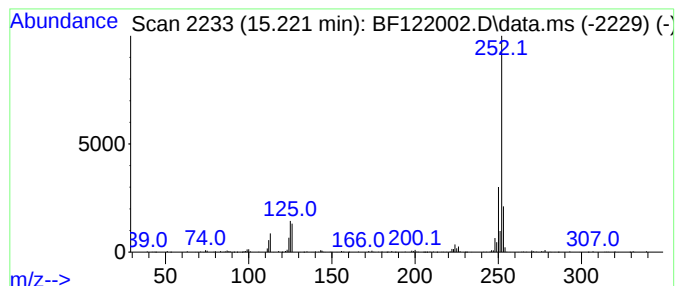
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Perylene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.221	7.98 ng	287306	Perylene-d12	15.339

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[e]pyrene	252	C20H12	000192-97-2	99
2			Benz[e]acephenanthrylene	252	C20H12	000205-99-2	98
3			Benzo[a]pyrene	252	C20H12	000050-32-8	96
4			Benzo[k]fluoranthene	252	C20H12	000207-08-9	94
5			Perylene	252	C20H12	000198-55-0	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF101520\
Data File : BF122002.D
Acq On : 15 Oct 2020 20:51
Operator : JU/CG
Sample : L4406-01 2X
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
JC-02-101420-A

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF101220.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Phenanthrene, 2...	11.792	2.4	ng	104158	4	11.263	875179	20.0
4H-Cyclopenta[d...	11.874	3.1	ng	136613	4	11.263	875179	20.0
9,12-Octadecadi...	12.333	4.3	ng	185877	4	11.263	875179	20.0
9-Octadecenoic ...	12.357	5.7	ng	248230	4	11.263	875179	20.0
Cyclopenta(def)...	12.404	2.0	ng	88810	4	11.263	875179	20.0
Fluoranthene, 2...	13.033	3.2	ng	203808	5	13.904	1277080	20.0
Benzo[e]pyrene	15.033	3.3	ng	117066	6	15.339	719816	20.0
Perylene	15.221	8.0	ng	287306	6	15.339	719816	20.0