

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF080221\
 Data File : BF124911.D
 Acq On : 2 Aug 2021 16:57
 Operator : JU/CG
 Sample : M3253-03
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MH-2-COMP

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-BF072221.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF124911.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.116	3	5	9	rVB	17315	16717	13.02%	1.148%
2	5.069	504	507	510	rBB	2293	2254	1.76%	0.155%
3	5.469	571	575	578	rBV	122459	102967	80.21%	7.070%
4	6.481	743	747	750	rBB	120675	106896	83.27%	7.340%
5	6.851	808	810	813	rBB	36017	27978	21.79%	1.921%
6	7.416	902	906	909	rBB	86076	71120	55.40%	4.884%
7	8.033	1009	1011	1014	rBB	2571	2184	1.70%	0.150%
8	8.133	1025	1028	1031	rBV	44122	38205	29.76%	2.623%
9	8.157	1031	1032	1034	rVB	3974	1951	1.52%	0.134%
10	8.845	1147	1149	1152	rBB	7708	6265	4.88%	0.430%
11	8.945	1164	1166	1169	rBB	14655	10626	8.28%	0.730%
12	9.210	1207	1211	1214	rBV	169405	128376	100.00%	8.815%
13	9.474	1253	1256	1259	rBB2	4095	4730	3.68%	0.325%
14	9.545	1265	1268	1271	rBV	13951	11683	9.10%	0.802%
15	9.574	1271	1273	1275	rVB	7898	5125	3.99%	0.352%
16	9.663	1286	1288	1293	rBB	6090	5235	4.08%	0.359%
17	9.745	1300	1302	1305	rBB2	7068	5634	4.39%	0.387%
18	9.886	1324	1326	1329	rVB	52487	43257	33.70%	2.970%
19	9.922	1330	1332	1334	rBB	2838	2023	1.58%	0.139%
20	10.086	1358	1360	1364	rBB2	4076	3474	2.71%	0.239%
21	10.127	1365	1367	1370	rBB	4196	3264	2.54%	0.224%
22	10.204	1377	1380	1382	rBB	3990	3389	2.64%	0.233%
23	10.233	1383	1385	1387	rBB	2947	1832	1.43%	0.126%
24	10.292	1392	1395	1397	rBV2	1822	2514	1.96%	0.173%
25	10.433	1416	1419	1423	rBB2	12791	11457	8.92%	0.787%
26	10.592	1444	1446	1448	rBB	2034	1685	1.31%	0.116%
27	10.674	1457	1460	1463	rBB	76225	55860	43.51%	3.836%
28	10.980	1508	1512	1515	rBB	6999	8454	6.59%	0.581%
29	11.010	1515	1517	1519	rBB	7603	4518	3.52%	0.310%
30	11.069	1524	1527	1529	rBB	3883	3329	2.59%	0.229%
31	11.180	1544	1546	1548	rBB	3629	2623	2.04%	0.180%
32	11.268	1559	1561	1564	rBB	6045	4306	3.35%	0.296%
33	11.374	1576	1579	1581	rBV	49494	40730	31.73%	2.797%
34	11.398	1581	1583	1587	rVB	80996	65202	50.79%	4.477%
35	11.451	1589	1592	1594	rBB	4008	3427	2.67%	0.235%
36	11.486	1595	1598	1600	rBB2	3102	3260	2.54%	0.224%

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 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

37	11.610	1617	1619	1623	rBB	1524	1648	1.28%	0.113%
38	11.704	1631	1635	1638	rBB	6913	5667	4.41%	0.389%
39	11.786	1646	1649	1652	rVB	4335	4055	3.16%	0.278%
40	11.874	1661	1664	1667	rBV	32312	25987	20.24%	1.784%
41	11.904	1667	1669	1672	rVB	38800	28151	21.93%	1.933%
42	11.951	1675	1677	1679	rBV2	4459	3326	2.59%	0.228%
43	11.986	1679	1683	1685	rVV2	30316	29668	23.11%	2.037%
44	12.010	1685	1687	1690	rVB	26605	20045	15.61%	1.376%
45	12.168	1712	1714	1716	rBV	8768	6067	4.73%	0.417%
46	12.198	1716	1719	1722	rVB2	5679	5669	4.42%	0.389%
47	12.321	1738	1740	1742	rVB	3967	2814	2.19%	0.193%
48	12.351	1742	1745	1747	rBV	7321	5775	4.50%	0.397%
49	12.374	1747	1749	1752	rVB2	3897	3020	2.35%	0.207%
50	12.427	1755	1758	1760	rBV	20350	17766	13.84%	1.220%
51	12.463	1760	1764	1766	rVV2	11220	14230	11.08%	0.977%
52	12.480	1766	1767	1770	rVV	7197	4256	3.32%	0.292%
53	12.510	1770	1772	1777	rVB2	6590	8875	6.91%	0.609%
54	12.586	1782	1785	1788	rVB	47381	36537	28.46%	2.509%
55	12.633	1789	1793	1794	rBV	2980	2063	1.61%	0.142%
56	12.651	1794	1796	1799	rVB2	4445	3208	2.50%	0.220%
57	12.686	1799	1802	1804	rBB	1974	1776	1.38%	0.122%
58	12.715	1805	1807	1809	rBB	2835	1822	1.42%	0.125%
59	12.739	1809	1811	1813	rBB	2169	1617	1.26%	0.111%
60	12.815	1819	1824	1830	rBB	57771	52632	41.00%	3.614%
61	12.868	1830	1833	1836	rBB2	3755	3869	3.01%	0.266%
62	12.927	1840	1843	1845	rBV2	1787	1818	1.42%	0.125%
63	12.957	1845	1848	1851	rBV	122988	95658	74.51%	6.569%
64	13.033	1857	1861	1868	rBB3	7017	9139	7.12%	0.628%
65	13.145	1873	1880	1883	rBB2	15095	20428	15.91%	1.403%
66	13.210	1889	1891	1894	rVB	8479	6432	5.01%	0.442%
67	13.245	1894	1897	1900	rBB2	13263	11136	8.67%	0.765%
68	13.339	1910	1913	1916	rBV	13145	10095	7.86%	0.693%
69	13.368	1916	1918	1922	rVB	10983	9212	7.18%	0.633%
70	13.545	1944	1948	1949	rBV	2411	2108	1.64%	0.145%
71	13.562	1949	1951	1955	rVB2	3834	3735	2.91%	0.256%
72	13.627	1959	1962	1964	rBV	2263	3358	2.62%	0.231%
73	13.651	1964	1966	1970	rVB3	5974	5982	4.66%	0.411%
74	13.710	1974	1976	1980	rBB	1955	1937	1.51%	0.133%
75	13.762	1980	1985	1988	rBV5	3859	6556	5.11%	0.450%
76	13.786	1988	1989	1992	rVB	3971	2388	1.86%	0.164%

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77	13.857	1999	2001	2005	rVB	2369	2656	2.07%	0.182%
78	14.009	2023	2027	2030	rBV2	37874	46467	36.20%	3.191%
79	14.039	2030	2032	2039	rVB	22562	18424	14.35%	1.265%
80	14.157	2050	2052	2053	rBV	2772	1728	1.35%	0.119%
81	14.398	2089	2093	2096	rBV	9761	9873	7.69%	0.678%
82	14.433	2096	2099	2101	rVB	5531	4762	3.71%	0.327%
83	14.468	2102	2105	2107	rBV	1923	1811	1.41%	0.124%
84	14.492	2107	2109	2111	rVB	3030	2227	1.73%	0.153%
85	14.521	2111	2114	2117	rBV	5789	5419	4.22%	0.372%
86	15.051	2199	2204	2207	rBV3	8804	12957	10.09%	0.890%
87	15.074	2207	2208	2211	rVB	4863	2836	2.21%	0.195%
88	15.156	2219	2222	2225	rBB	1962	1635	1.27%	0.112%
89	15.351	2252	2255	2259	rBB	7984	7438	5.79%	0.511%
90	15.415	2262	2266	2270	rBB	10877	11175	8.70%	0.767%
91	15.480	2273	2277	2281	rBV	22058	25994	20.25%	1.785%
92	16.939	2520	2525	2530	rBB2	2616	4207	3.28%	0.289%
93	17.386	2596	2601	2604	rBB	2704	3612	2.81%	0.248%

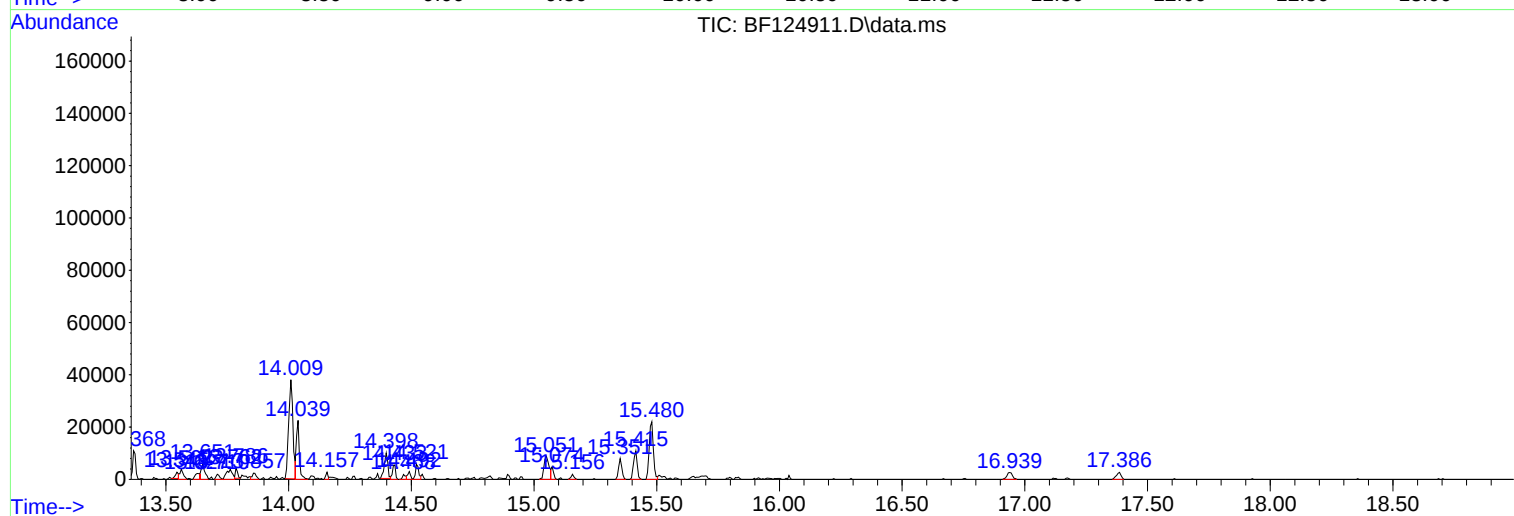
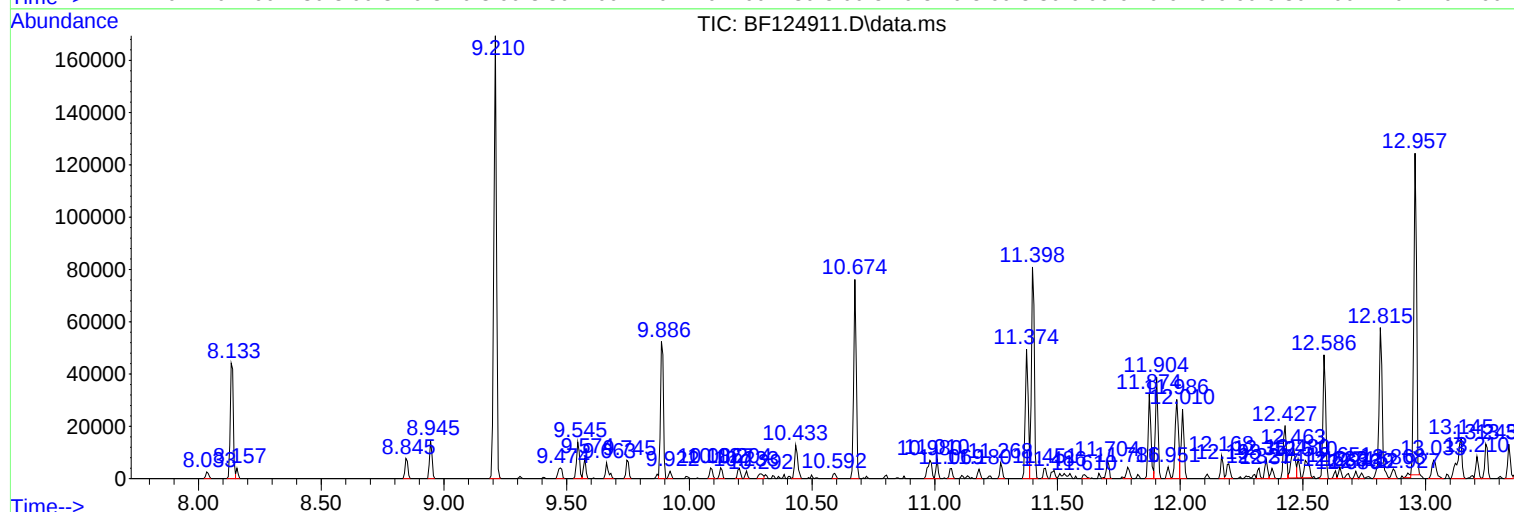
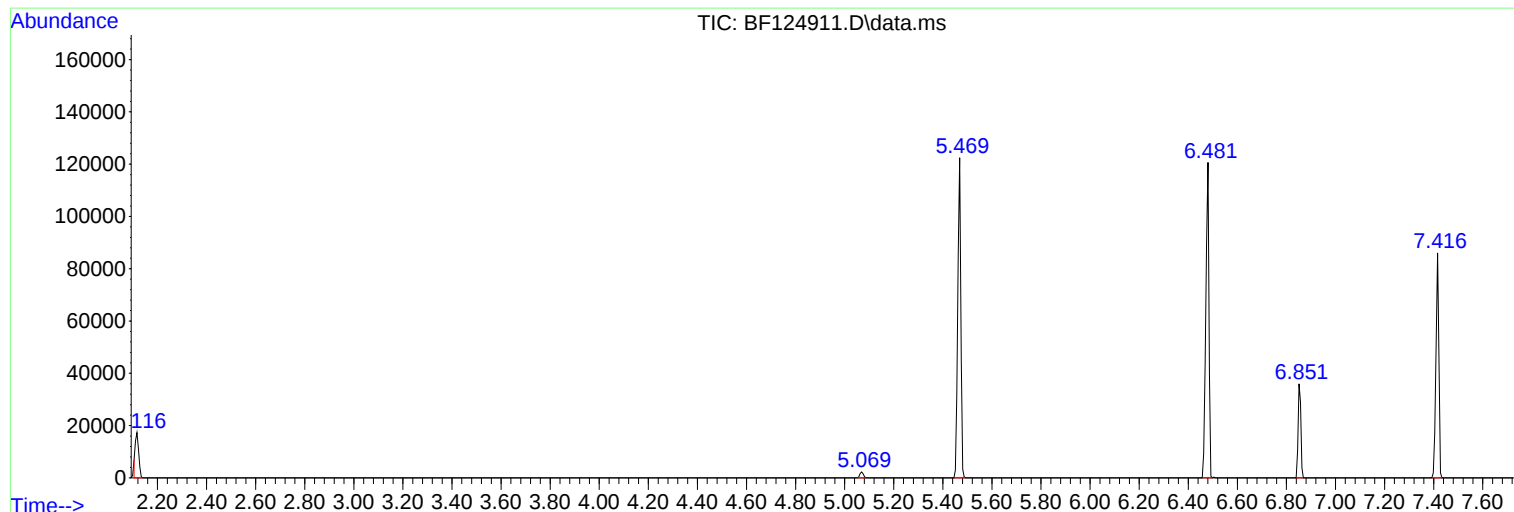
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BNA_F
ClientSampled :
MH-2-COMP

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF072221.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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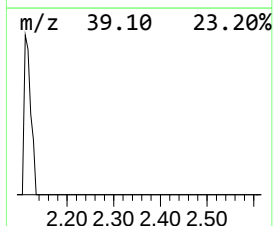
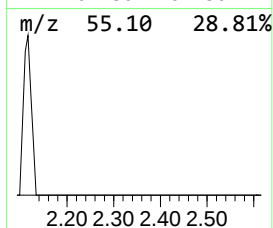
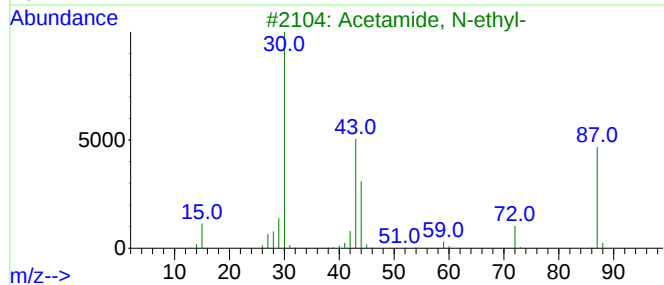
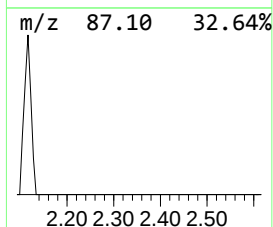
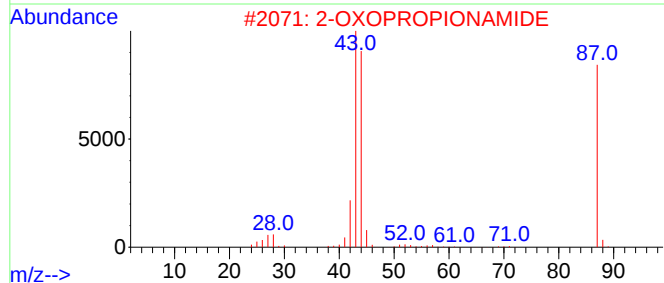
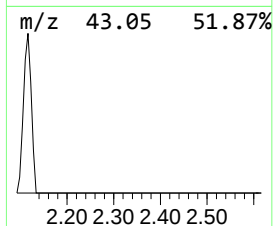
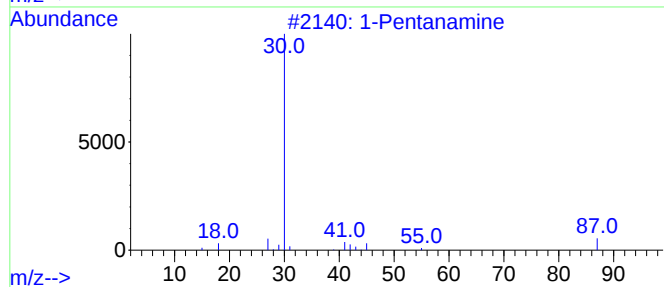
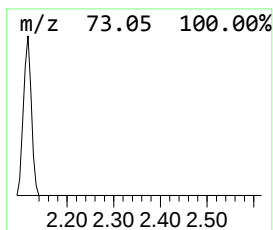
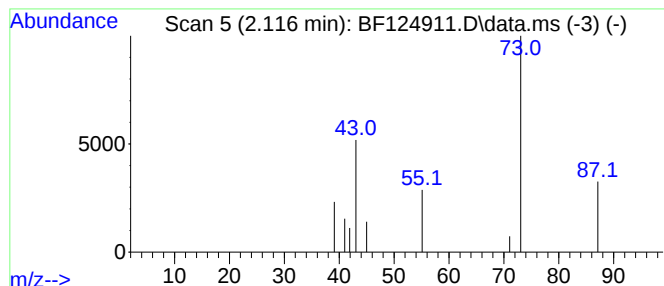
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF072221.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 unknown2.116 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.116	11.95 ng	16717	1,4-Dichlorobenzene-d4	6.851

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Pentanamine	87	C5H13N	000110-58-7	5
2			2-OXOPROPIONAMIDE	87	C3H5NO2	000631-66-3	5
3			Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	5
4			Acetamide, N-methyl-	73	C3H7NO	000079-16-3	5
5			Isobutylamine	73	C4H11N	000078-81-9	5



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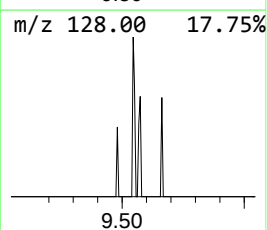
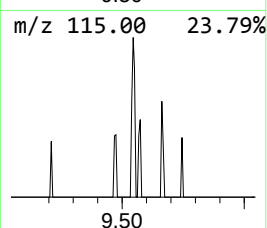
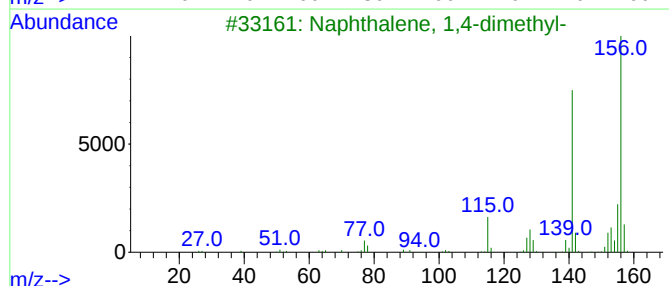
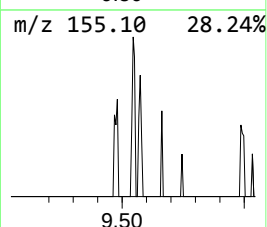
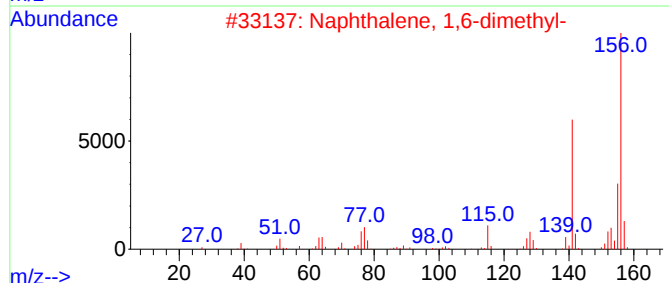
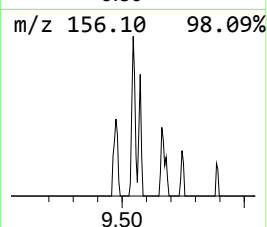
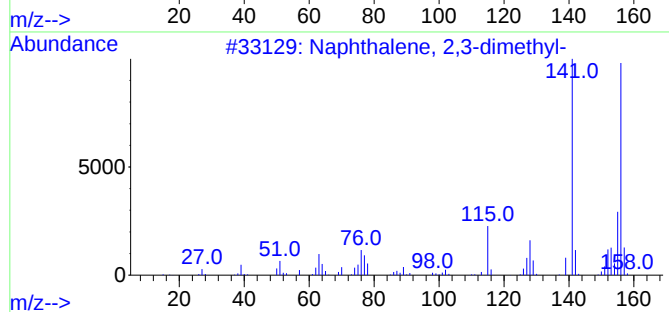
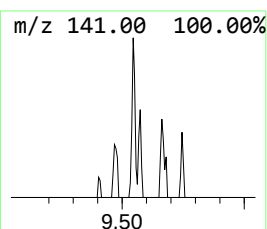
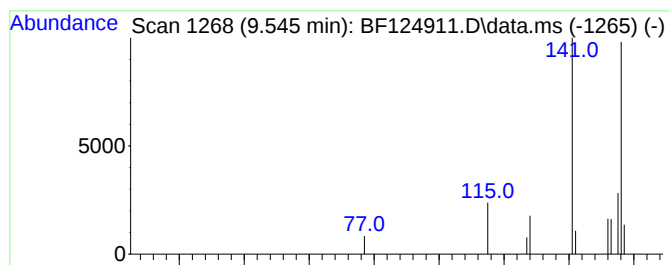
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF072221.M
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Naphthalene, 2,3-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.545	5.40 ng	11683	Acenaphthene-d10	9.892

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	94
2			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	94
3			Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	91
4			Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	91
5			Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	91



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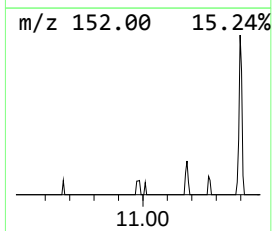
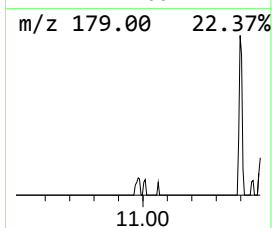
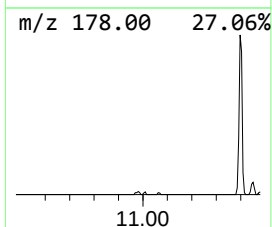
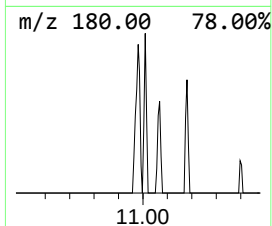
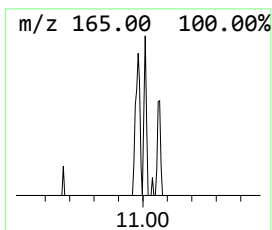
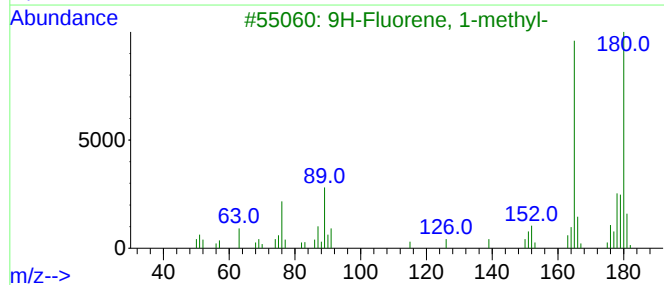
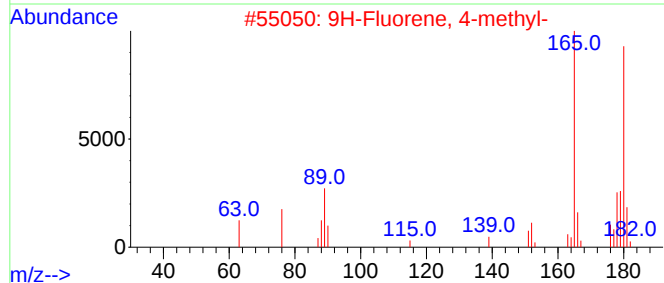
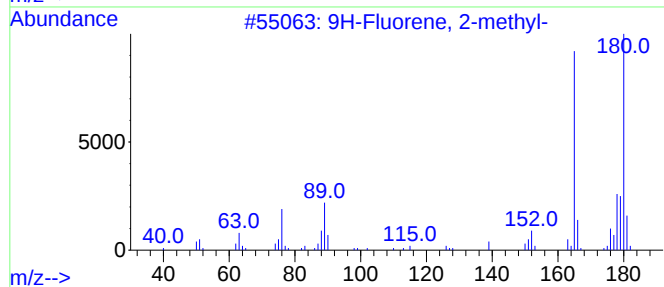
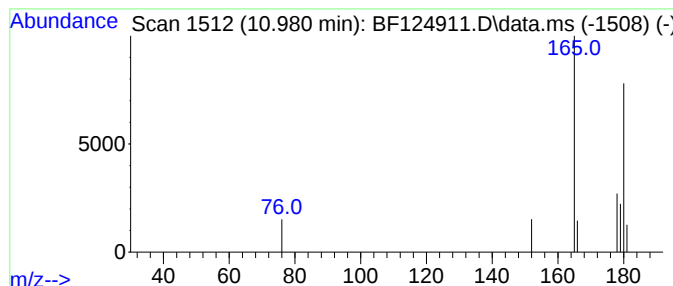
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ClientSampleId :
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 9H-Fluorene, 2-methyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.980	4.15 ng	8454	Phenanthrene-d10	11.374	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9H-Fluorene, 2-methyl-	180	C14H12	001430-97-3	87
2	9H-Fluorene, 4-methyl-	180	C14H12	001556-99-6	87
3	9H-Fluorene, 1-methyl-	180	C14H12	001730-37-6	87
4	9H-Fluorene, 9-methyl-	180	C14H12	002523-37-7	86
5	6-Fluoro[1,2,4]triazolo[1,5-a]py...	180	C8H9FN4	1641556-52-6	86



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Data File : BF124911.D
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Operator : JU/CG
Sample : M3253-03
Misc :
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Instrument :
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ClientSampleId :
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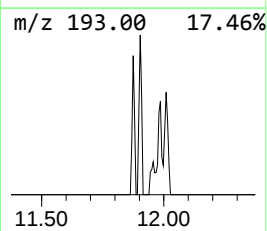
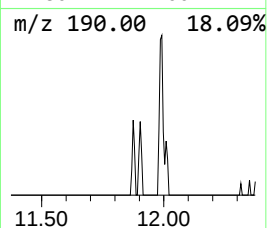
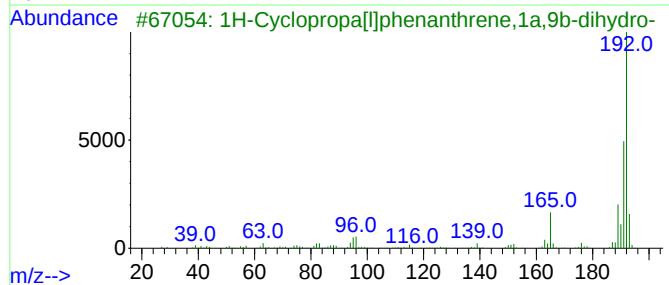
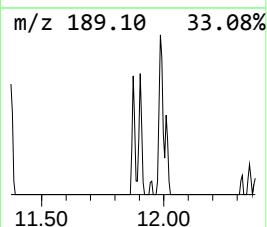
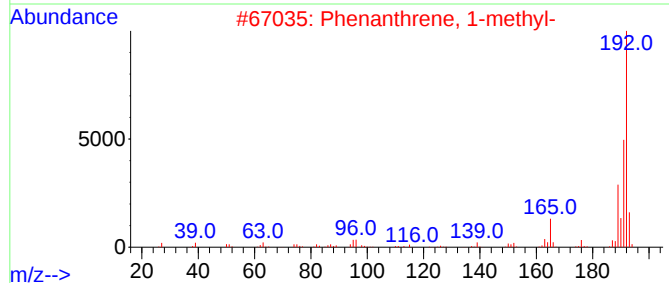
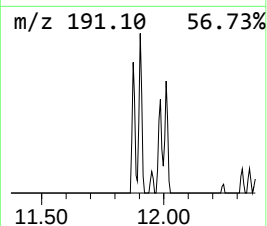
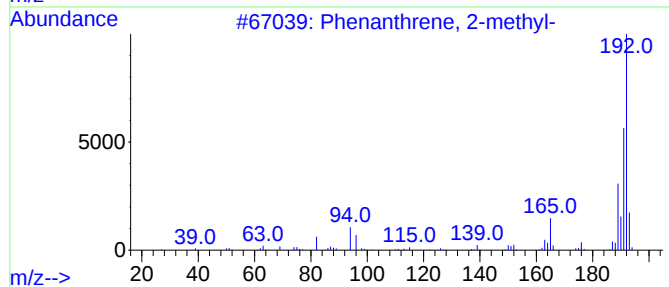
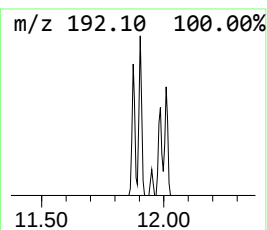
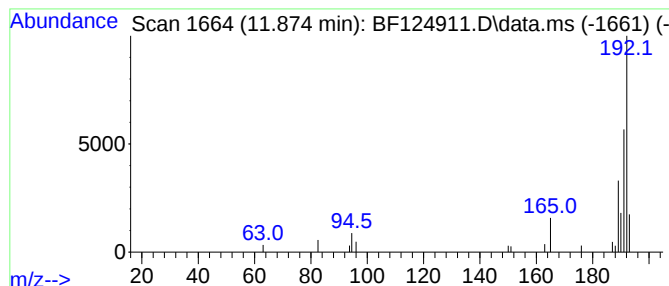
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 4 Phenanthrene, 1-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.874	12.76 ng	25987	Phenanthrene-d10	11.374

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	90
2			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	87
3			1H-Cyclopropa[l]phenanthrene,1a,...	192	C15H12	000949-41-7	87
4			Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	83
5			Anthracene, 1-methyl-	192	C15H12	000610-48-0	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF080221\
Data File : BF124911.D
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Sample : M3253-03
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ClientSampleId :
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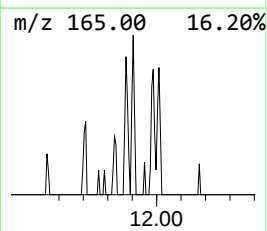
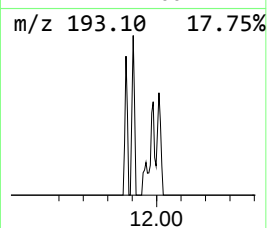
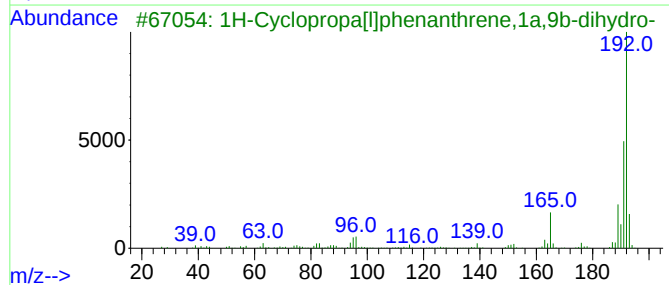
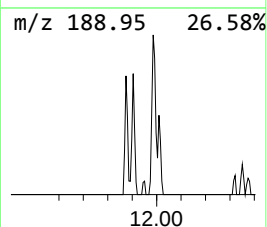
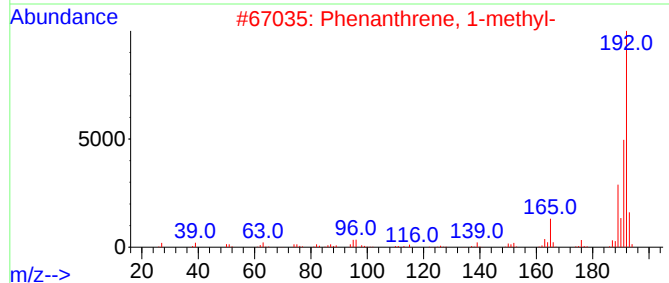
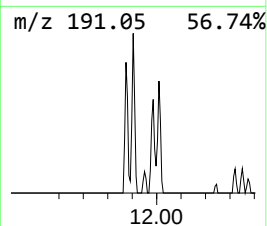
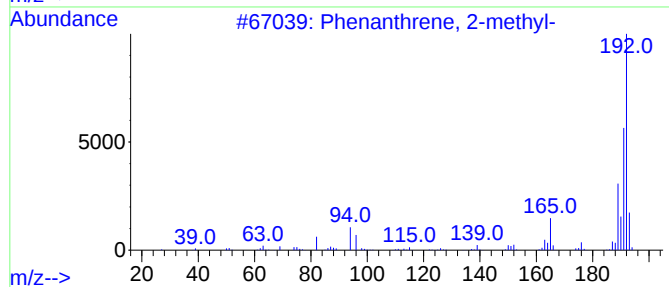
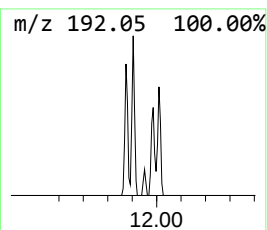
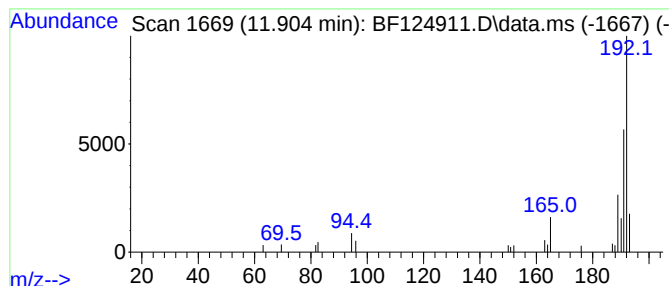
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.904	13.82 ng	28151	Phenanthrene-d10	11.374

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	96
2			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	95
3			1H-Cyclopropa[1]phenanthrene,1a,...	192	C15H12	000949-41-7	94
4			Anthracene, 1-methyl-	192	C15H12	000610-48-0	94
5			1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF080221\
Data File : BF124911.D
Acq On : 2 Aug 2021 16:57
Operator : JU/CG
Sample : M3253-03
Misc :
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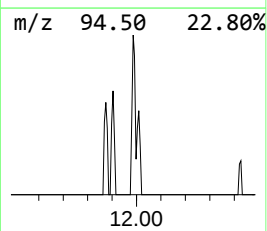
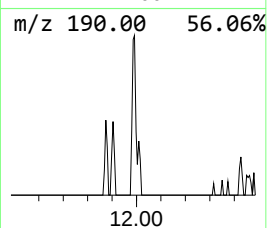
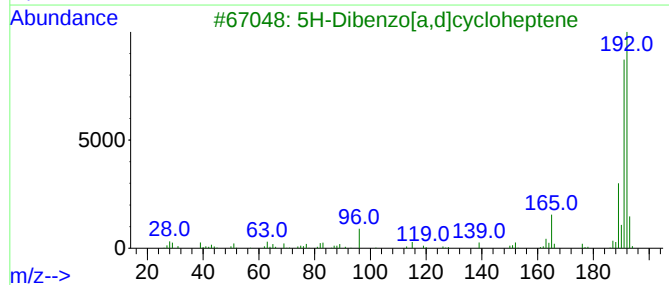
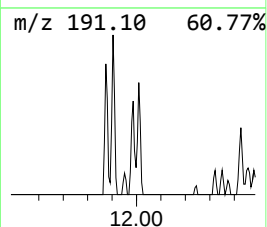
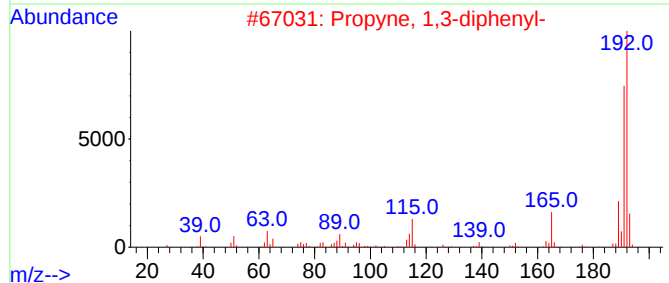
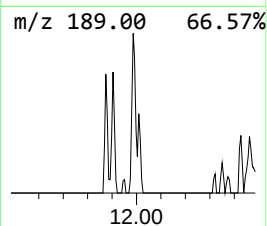
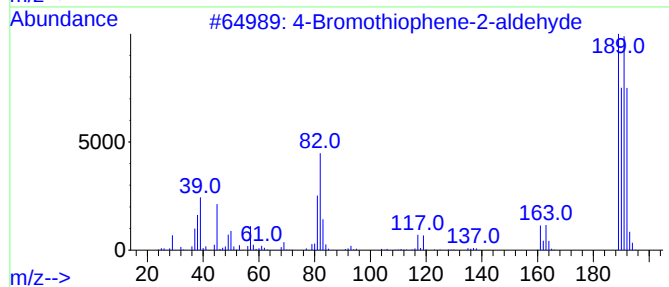
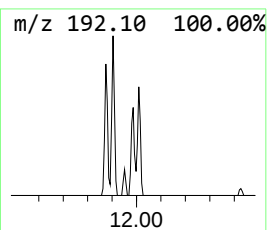
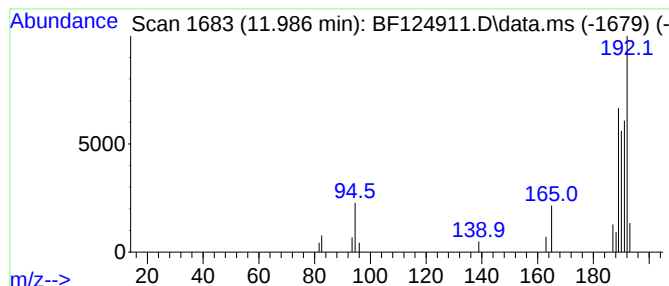
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF072221.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 unknown11.986 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.986	14.57 ng	29668	Phenanthrene-d10	11.374	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-Bromothiophene-2-aldehyde	190	C5H3BrOS	018791-75-8	49
2	Propyne, 1,3-diphenyl-	192	C15H12	004980-70-5	49
3	5H-Dibenzo[a,d]cycloheptene	192	C15H12	000256-81-5	47
4	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	47
5	1H-Cyclopropa[1]phenanthrene,1a,...	192	C15H12	000949-41-7	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF080221\
Data File : BF124911.D
Acq On : 2 Aug 2021 16:57
Operator : JU/CG
Sample : M3253-03
Misc :
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Instrument :
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ClientSampleId :
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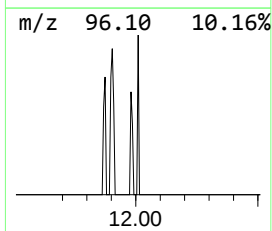
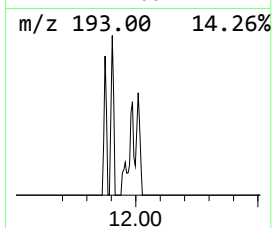
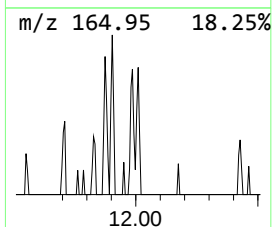
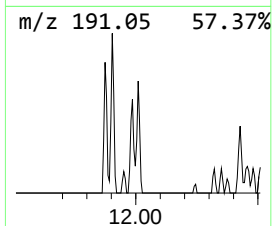
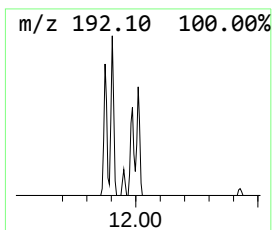
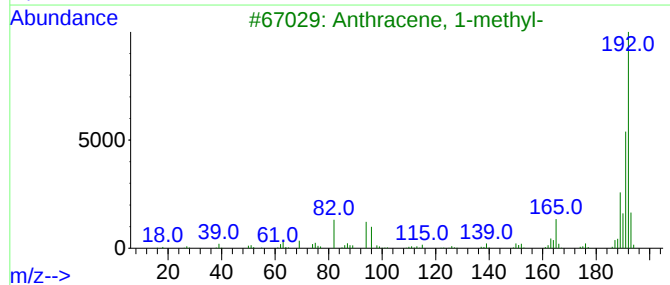
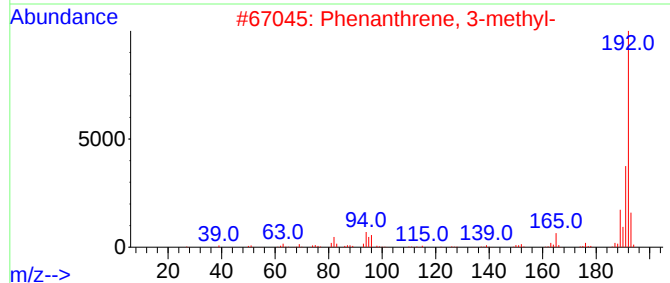
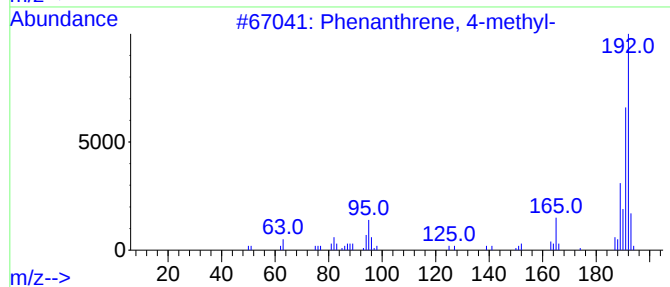
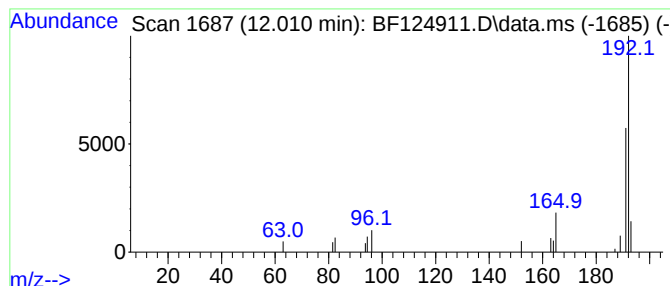
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Phenanthrene, 4-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.010	9.84 ng	20045	Phenanthrene-d10	11.374

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	91
2			Phenanthrene, 3-methyl-	192	C15H12	000832-71-3	91
3			Anthracene, 1-methyl-	192	C15H12	000610-48-0	91
4			Anthracene, 9-methyl-	192	C15H12	000779-02-2	90
5			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF080221\
Data File : BF124911.D
Acq On : 2 Aug 2021 16:57
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Misc :
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Instrument :
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ClientSampleId :
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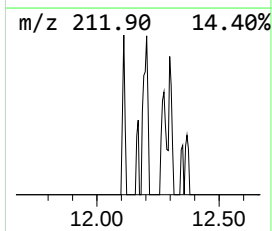
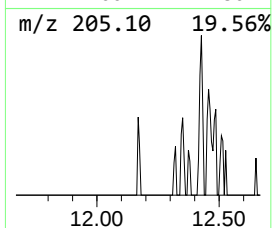
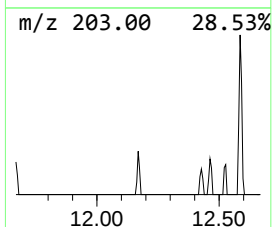
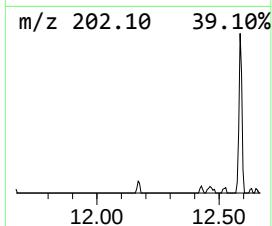
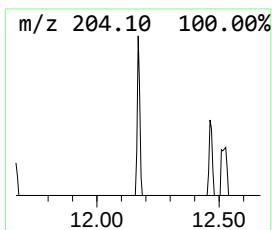
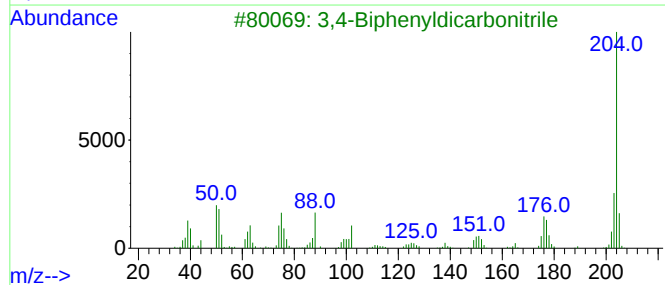
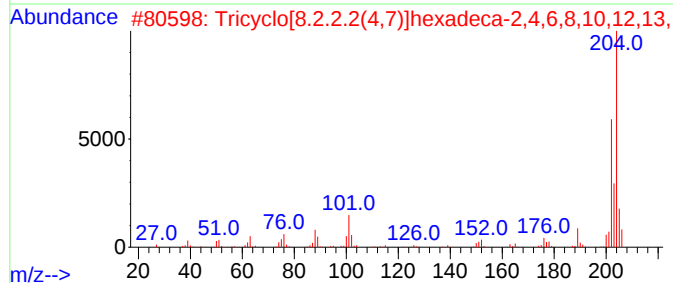
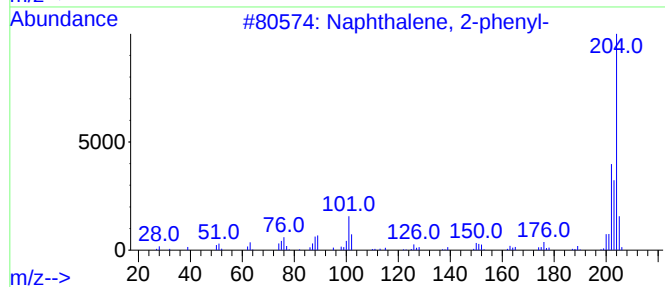
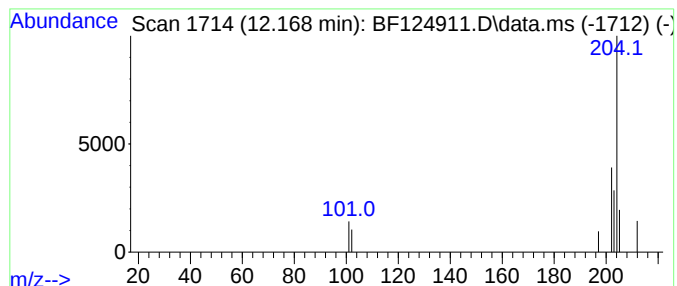
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Naphthalene, 2-phenyl- Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.168	2.98 ng	6067	Phenanthrene-d10	11.374

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	58
2			Tricyclo[8.2.2.2(4,7)]hexadeca-2...	204	C16H12	006572-60-7	53
3			3,4-Biphenyldicarbonitrile	204	C14H8N2	004128-63-6	53
4			5-Amino-6,8-dimethoxyquinoline	204	C11H12N2O2	1000213-95-2	53
5			[1,1'-Biphenyl]-2-ol, 3-chloro-	204	C12H9ClO	000085-97-2	42



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF080221\
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ClientSampleId :
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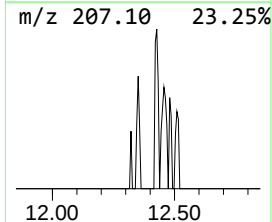
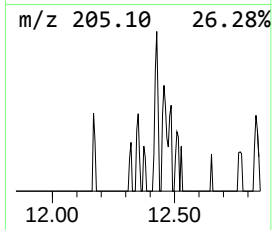
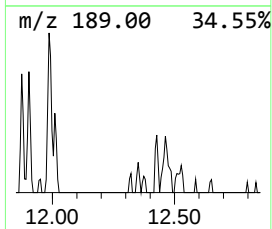
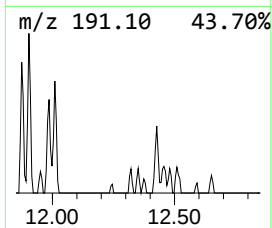
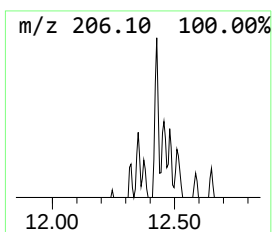
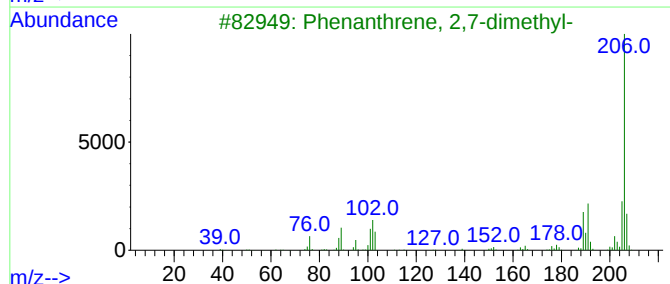
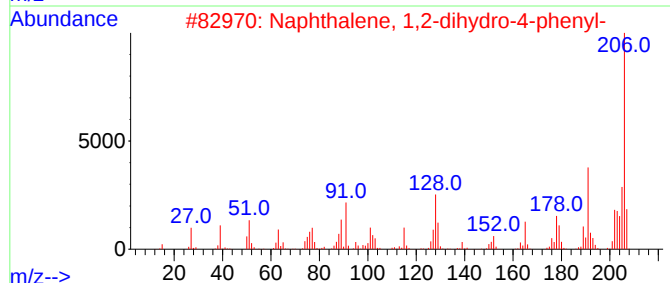
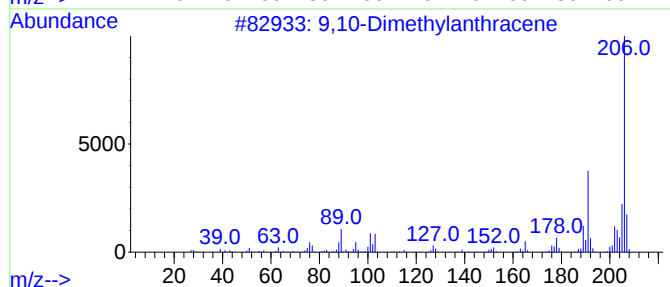
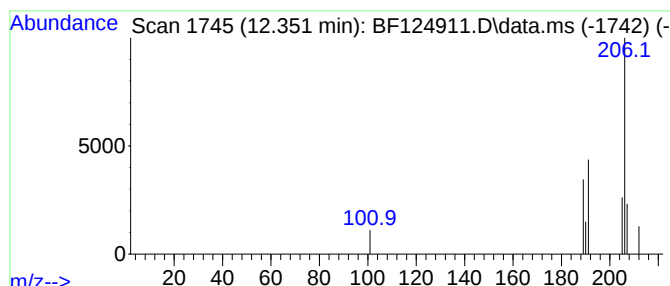
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 9,10-Dimethylantracene Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.351	2.84 ng	5775	Phenanthrene-d10	11.374

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		9,10-Dimethylantracene	206	C16H14	000781-43-1	90
2		Naphthalene, 1,2-dihydro-4-phenyl-	206	C16H14	007469-40-1	90
3		Phenanthrene, 2,7-dimethyl-	206	C16H14	001576-69-8	83
4		Phenanthrene, 1,7-dimethyl-	206	C16H14	000483-87-4	83
5		Phenanthrene, 3,6-dimethyl-	206	C16H14	001576-67-6	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF080221\
Data File : BF124911.D
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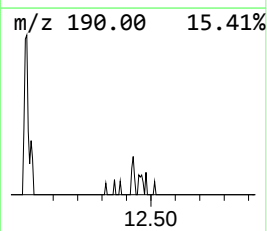
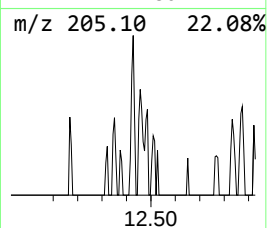
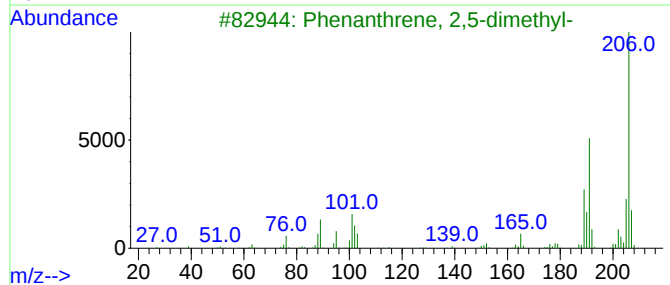
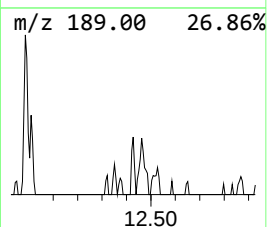
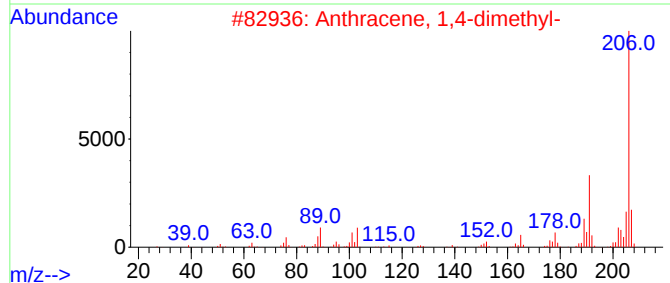
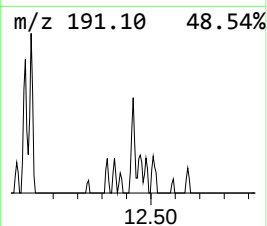
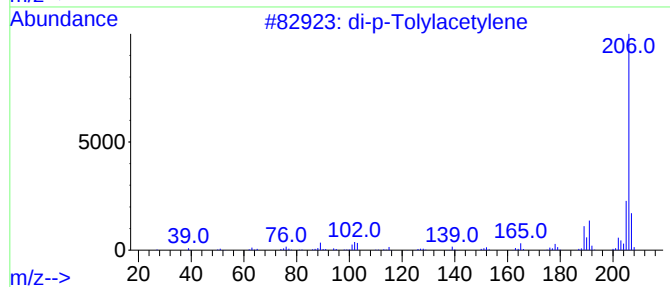
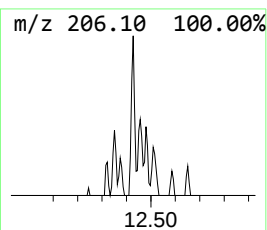
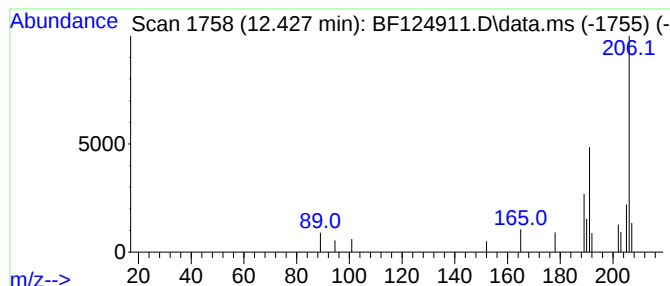
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 di-p-Tolylacetylene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.427	8.72 ng	17766	Phenanthrene-d10	11.374

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			di-p-Tolylacetylene	206	C16H14	002789-88-0	90
2			Anthracene, 1,4-dimethyl-	206	C16H14	000781-92-0	87
3			Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	87
4			9,10-Dimethylanthracene	206	C16H14	000781-43-1	87
5			3,4-Dimethylphenanthrene	206	C16H14	1000452-24-5	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF080221\
Data File : BF124911.D
Acq On : 2 Aug 2021 16:57
Operator : JU/CG
Sample : M3253-03
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MH-2-COMP

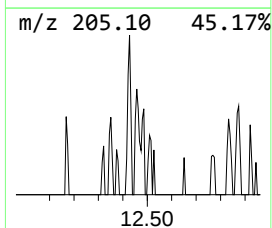
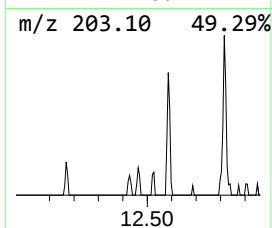
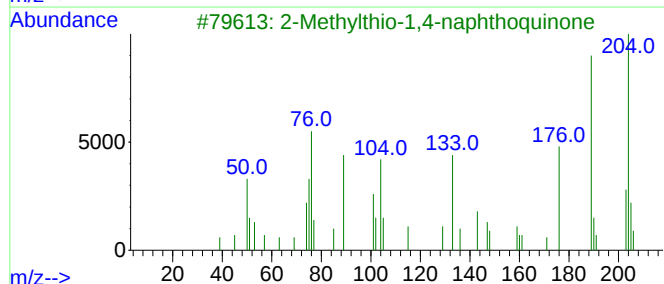
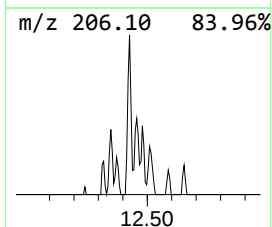
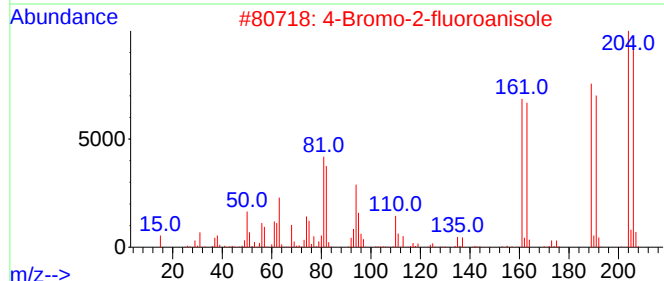
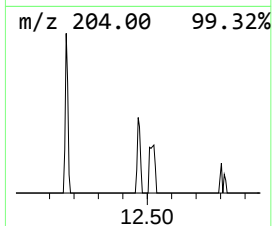
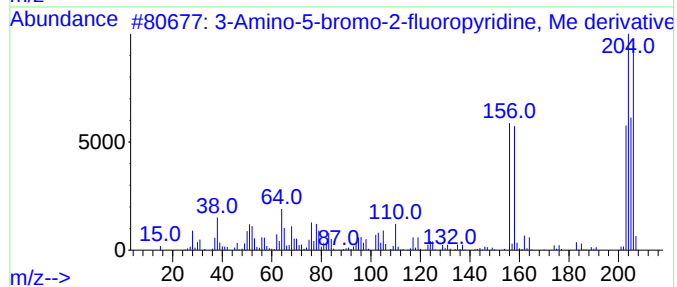
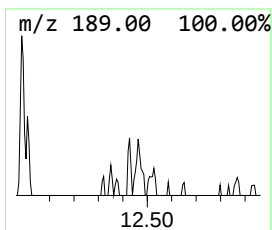
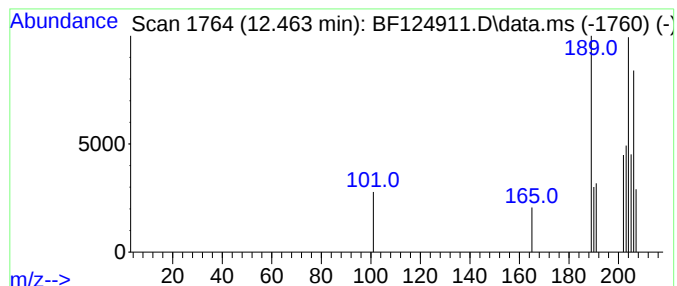
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 3-Amino-5-bromo-2-fluoropyr... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.463	6.99 ng	14230	Phenanthrene-d10	11.374

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Amino-5-bromo-2-fluoropyridine...	204	C6H6BrFN2	1000496-58-8	50
2			4-Bromo-2-fluoroanisole	204	C7H6BrFO	002357-52-0	33
3			2-Methylthio-1,4-naphthoquinone	204	C11H8O2S	026037-60-5	32
4			1,4-Diphenyl-1,3-butadiene	206	C16H14	000886-65-7	30
5			Quinoline, 8-methoxy-5-nitro-	204	C10H8N2O3	1000298-88-7	27



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF080221\
Data File : BF124911.D
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Misc :
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ClientSampleId :
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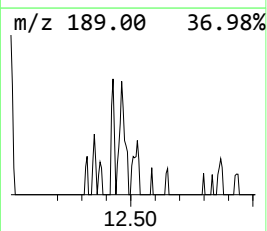
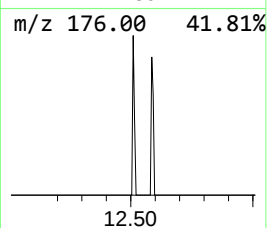
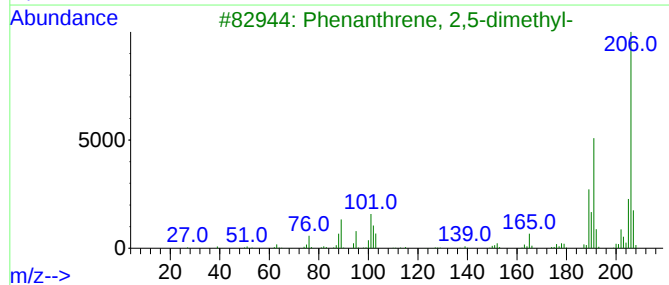
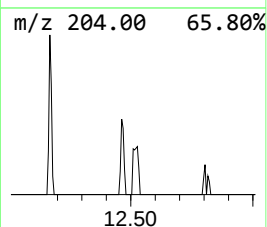
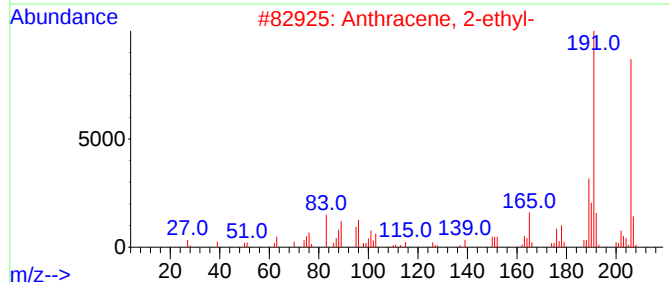
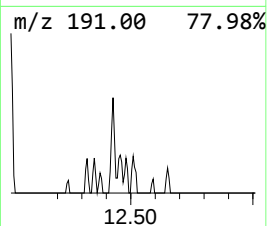
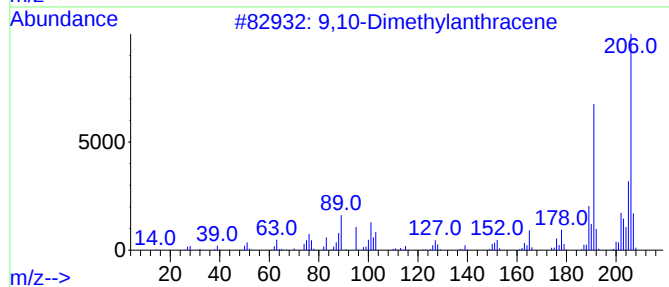
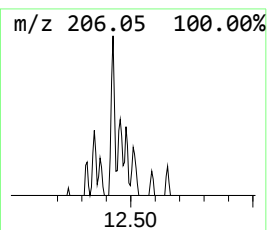
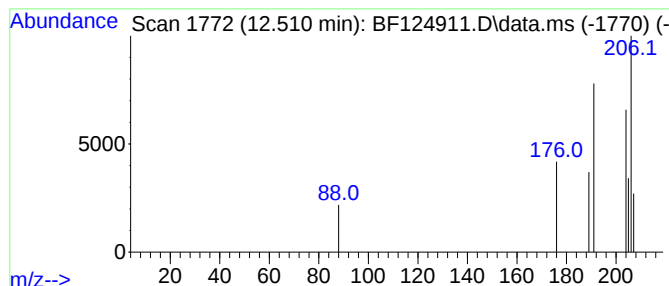
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 unknown12.510 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.510	4.36 ng	8875	Phenanthrene-d10	11.374

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		9,10-Dimethylantracene	206	C16H14	000781-43-1	42
2		Anthracene, 2-ethyl-	206	C16H14	052251-71-5	40
3		Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	40
4		Naphthalene, 1,2-dihydro-4-phenyl-	206	C16H14	007469-40-1	40
5		Phenanthrene-9-carboxaldehyde	206	C15H10O	004707-71-5	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF080221\
Data File : BF124911.D
Acq On : 2 Aug 2021 16:57
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Sample : M3253-03
Misc :
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Instrument :
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ClientSampleId :
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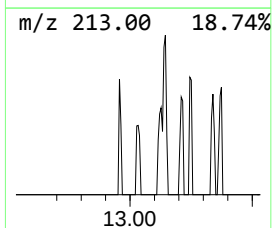
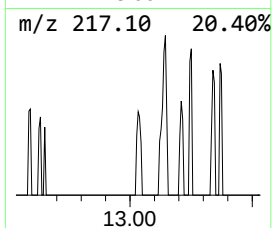
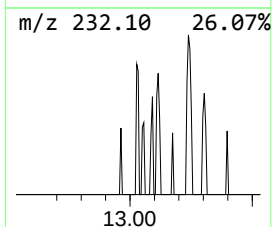
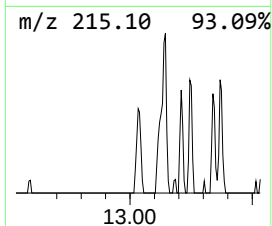
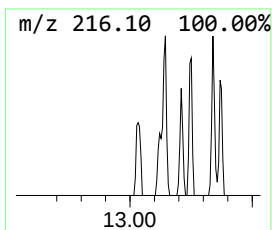
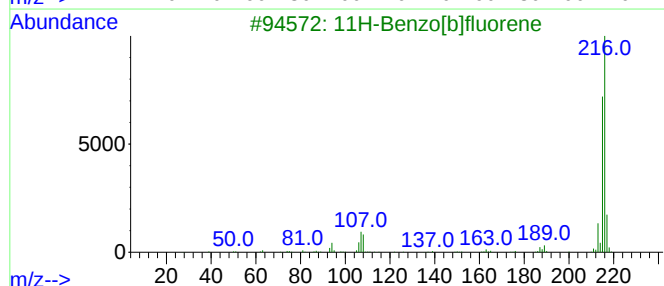
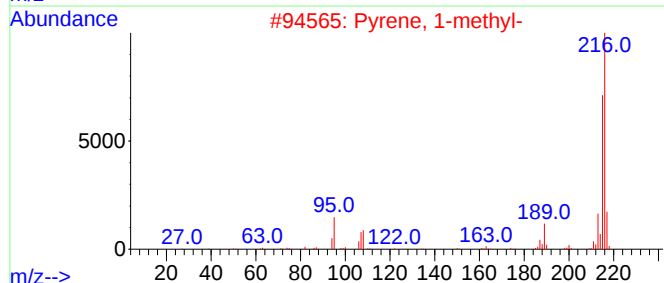
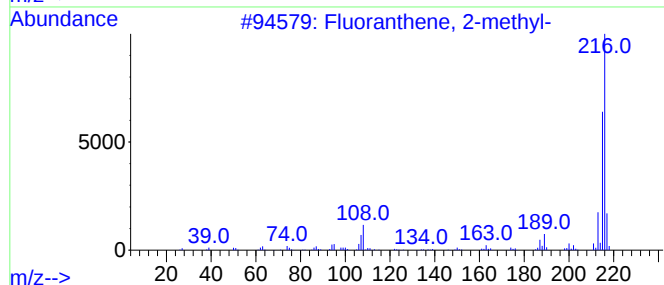
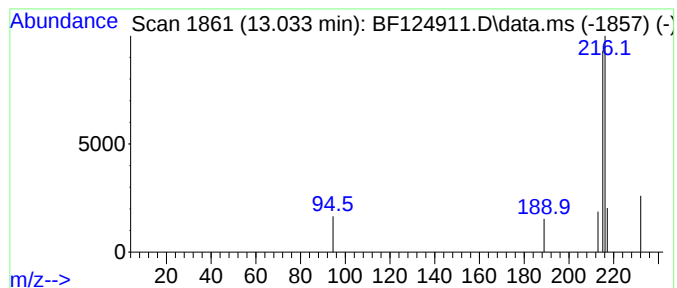
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Fluoranthene, 2-methyl- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.033	3.93 ng	9139	Chrysene-d12	14.009

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	86
2			Pyrene, 1-methyl-	216	C17H12	002381-21-7	86
3			11H-Benzo[b]fluorene	216	C17H12	000243-17-4	86
4			11H-Benzo[a]fluorene	216	C17H12	000238-84-6	78
5			1,3,5-Triazine-2,4,6-triamine, N...	216	C10H12N6	046731-79-7	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF080221\
Data File : BF124911.D
Acq On : 2 Aug 2021 16:57
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Misc :
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Instrument :
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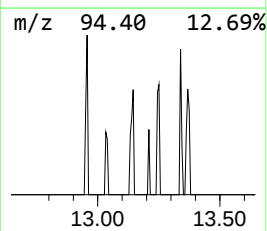
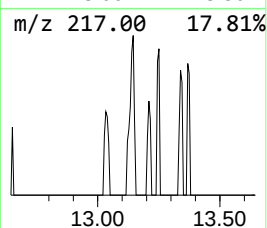
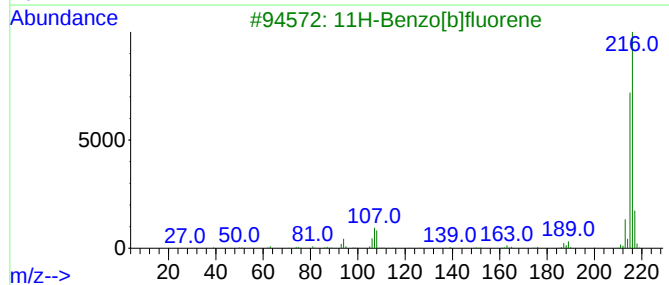
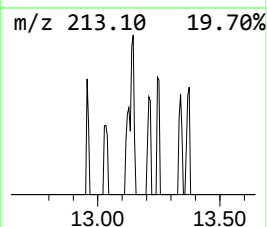
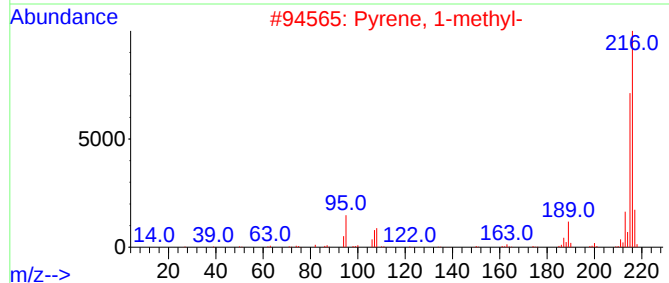
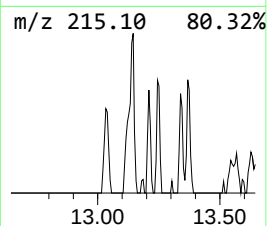
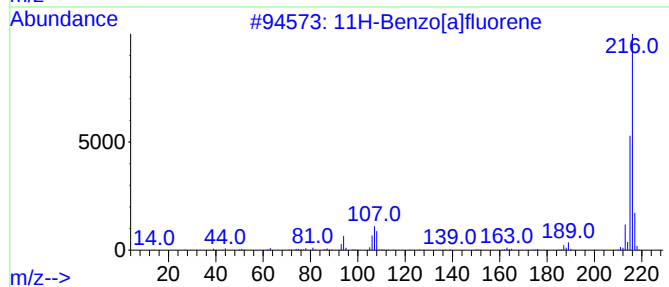
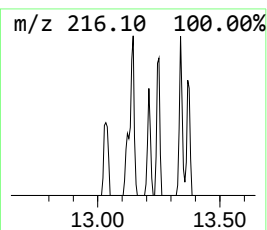
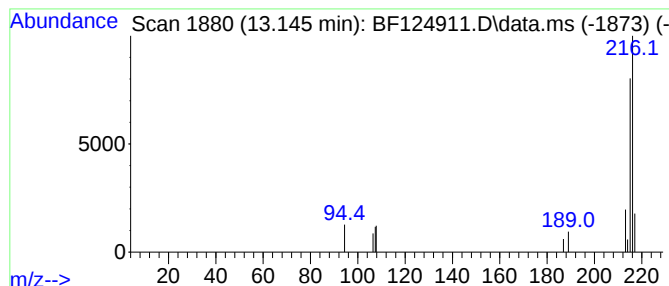
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 11H-Benzo[a]fluorene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.145	8.79 ng	20428	Chrysene-d12	14.009

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			11H-Benzo[a]fluorene	216	C17H12	000238-84-6	91
2			Pyrene, 1-methyl-	216	C17H12	002381-21-7	91
3			11H-Benzo[b]fluorene	216	C17H12	000243-17-4	91
4			Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	87
5			7H-Benzo[c]fluorene	216	C17H12	000205-12-9	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF080221\
Data File : BF124911.D
Acq On : 2 Aug 2021 16:57
Operator : JU/CG
Sample : M3253-03
Misc :
ALS Vial : 15 Sample Multiplier: 1

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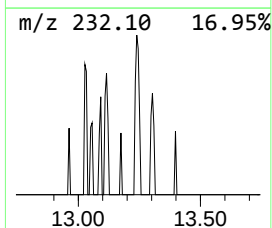
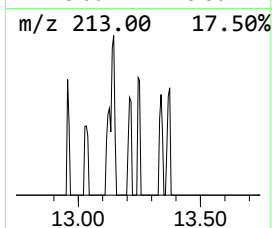
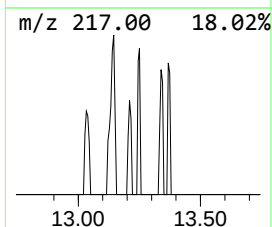
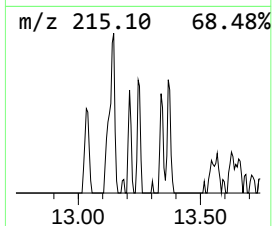
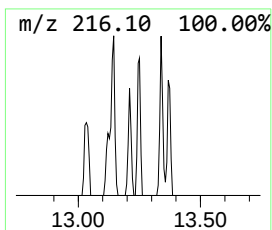
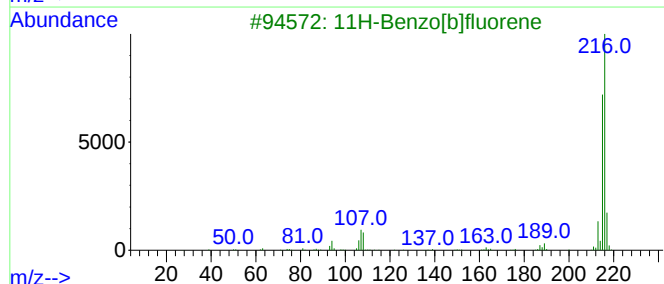
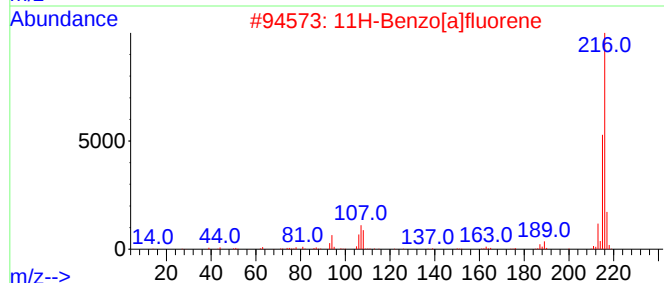
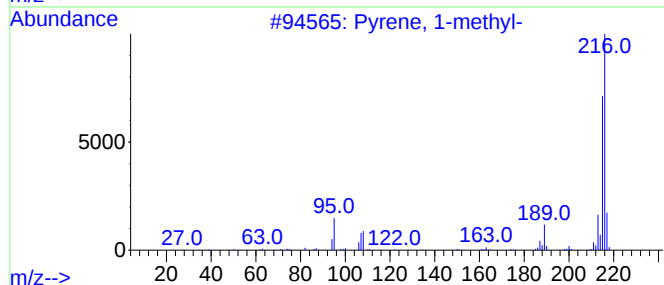
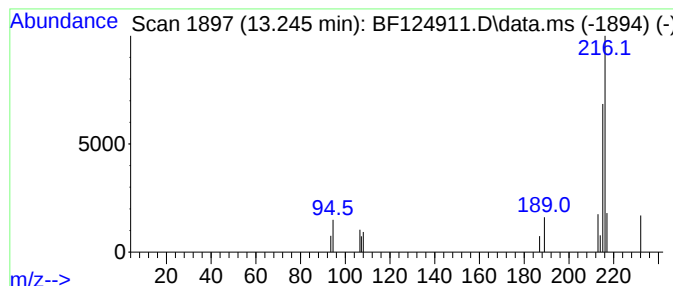
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Pyrene, 1-methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.245	4.79 ng	11136	Chrysene-d12	14.009

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pyrene, 1-methyl-	216	C17H12	002381-21-7	91
2			11H-Benzo[a]fluorene	216	C17H12	000238-84-6	86
3			11H-Benzo[b]fluorene	216	C17H12	000243-17-4	86
4			Pyrene, 2-methyl-	216	C17H12	003442-78-2	80
5			Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF080221\
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Misc :
ALS Vial : 15 Sample Multiplier: 1

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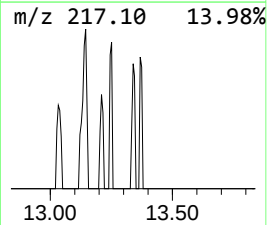
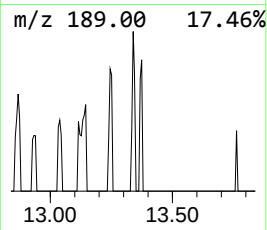
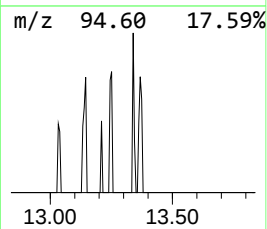
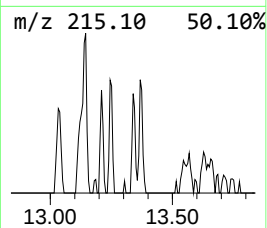
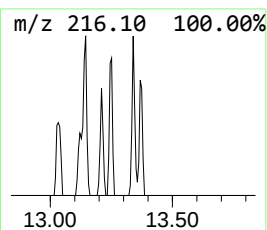
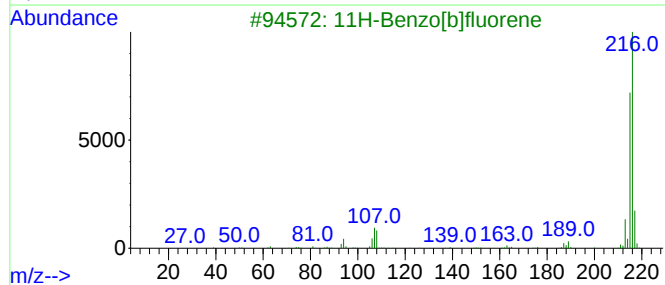
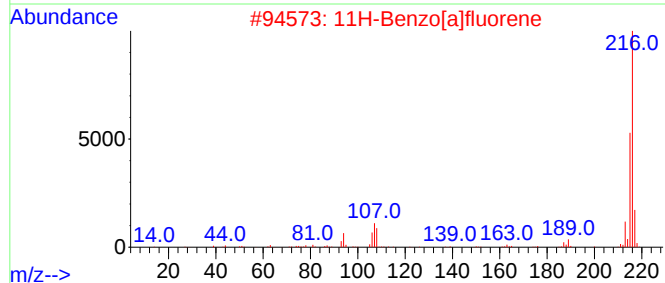
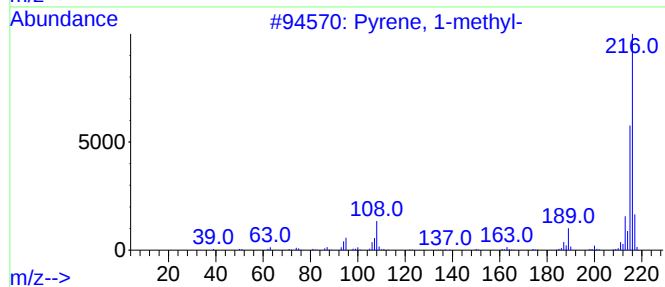
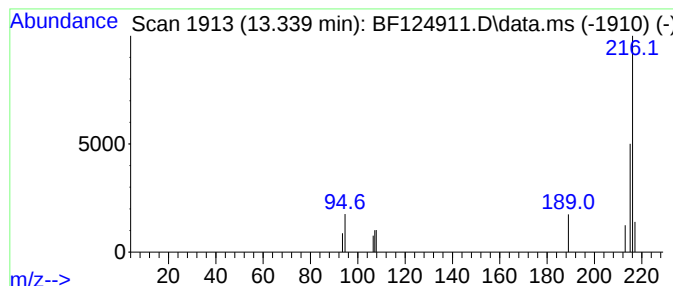
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 Pyrene, 4-methyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.339	4.35 ng	10095	Chrysene-d12	14.009

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Pyrene, 1-methyl-	216	C17H12	002381-21-7	90
2		11H-Benzo[a]fluorene	216	C17H12	000238-84-6	90
3		11H-Benzo[b]fluorene	216	C17H12	000243-17-4	90
4		Pyrene, 2-methyl-	216	C17H12	003442-78-2	83
5		Pyrene, 4-methyl-	216	C17H12	003353-12-6	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF080221\
Data File : BF124911.D
Acq On : 2 Aug 2021 16:57
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Sample : M3253-03
Misc :
ALS Vial : 15 Sample Multiplier: 1

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ClientSampleId :
MH-2-COMP

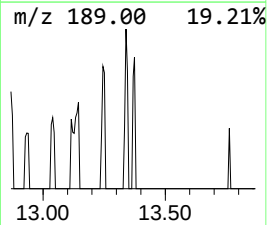
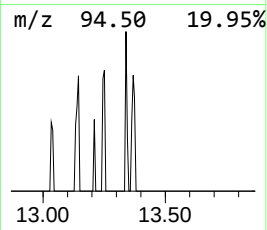
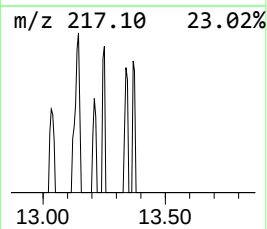
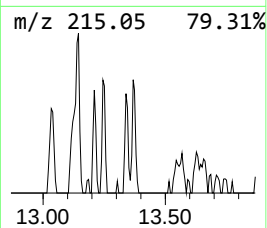
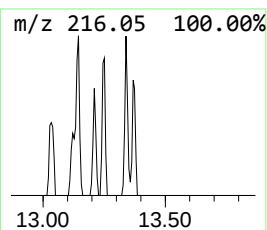
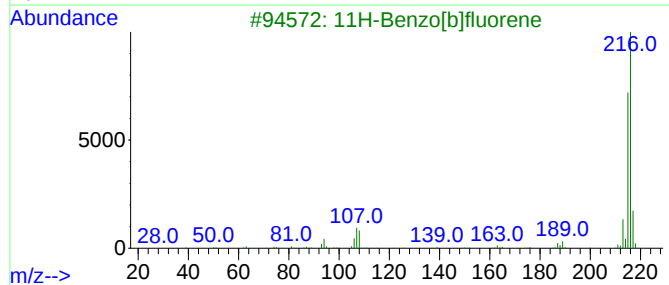
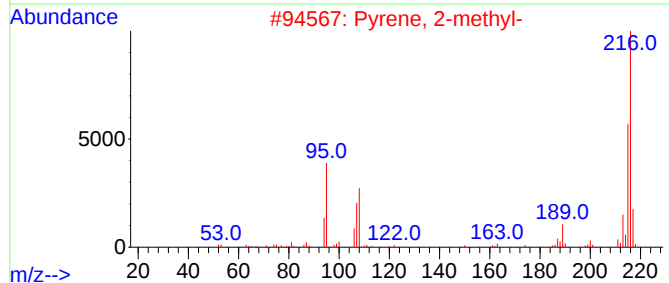
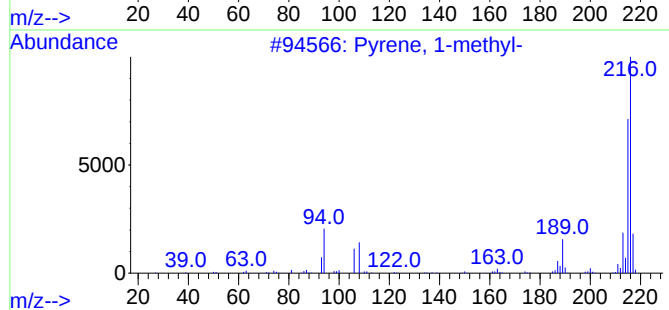
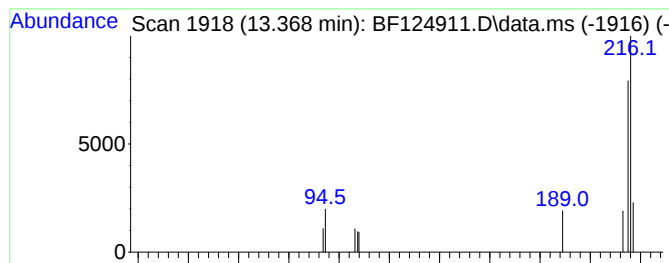
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 11H-Benzo[b]fluorene Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.368	3.96 ng	9212	Chrysene-d12	14.009

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pyrene, 1-methyl-		216	C17H12		002381-21-7	86
2	Pyrene, 2-methyl-		216	C17H12		003442-78-2	83
3	11H-Benzo[b]fluorene		216	C17H12		000243-17-4	80
4	Fluoranthene, 2-methyl-		216	C17H12		033543-31-6	59
5	11H-Benzo[a]fluorene		216	C17H12		000238-84-6	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF080221\
Data File : BF124911.D
Acq On : 2 Aug 2021 16:57
Operator : JU/CG
Sample : M3253-03
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MH-2-COMP

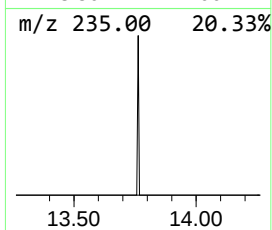
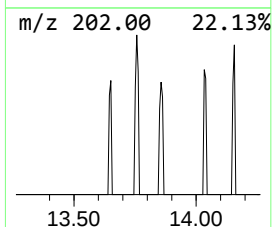
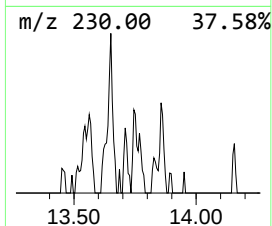
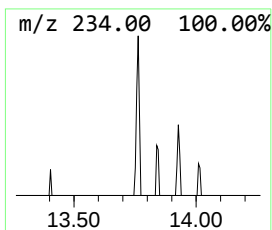
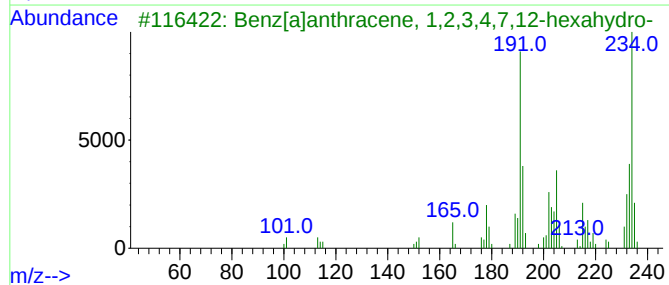
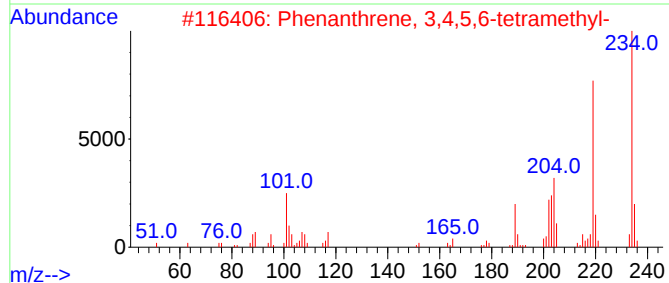
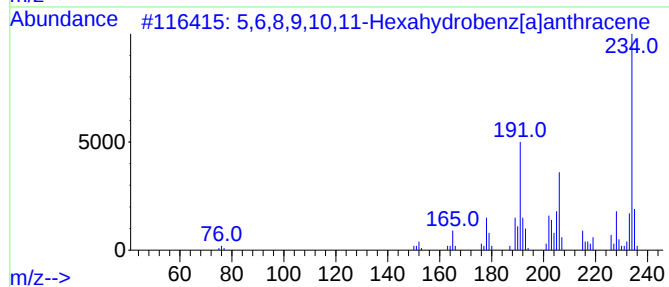
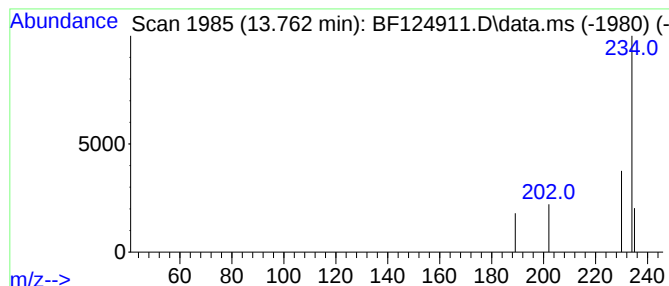
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 5,6,8,9,10,11-Hexahydrobenz... Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.762	2.82 ng	6556	Chrysene-d12	14.009

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	5,6,8,9,10,11-Hexahydrobenz[a]an...	234	C18H18	067064-61-3	72		
2	Phenanthrene, 3,4,5,6-tetramethyl-	234	C18H18	007343-06-8	64		
3	Benz[a]anthracene, 1,2,3,4,7,12-...	234	C18H18	016434-62-1	64		
4	Phenanthrene, 2,4,5,7-tetramethyl-	234	C18H18	007396-38-5	64		
5	Pyreno[4,5-b:9,10-b']bisoxirene,...	234	C16H10O2	055400-88-9	50		



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF080221\
Data File : BF124911.D
Acq On : 2 Aug 2021 16:57
Operator : JU/CG
Sample : M3253-03
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MH-2-COMP

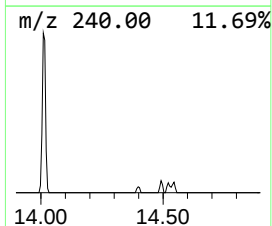
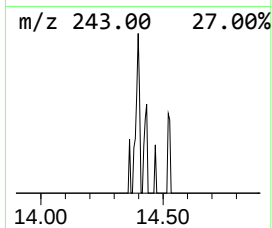
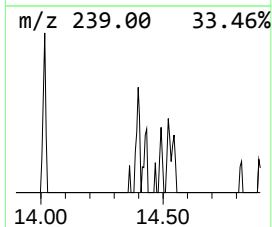
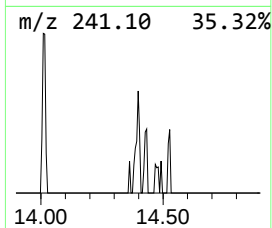
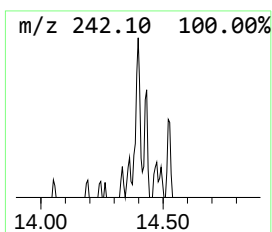
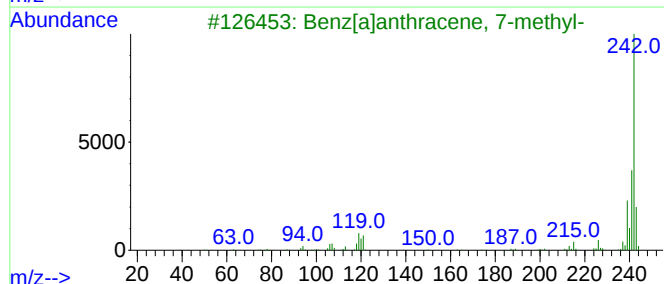
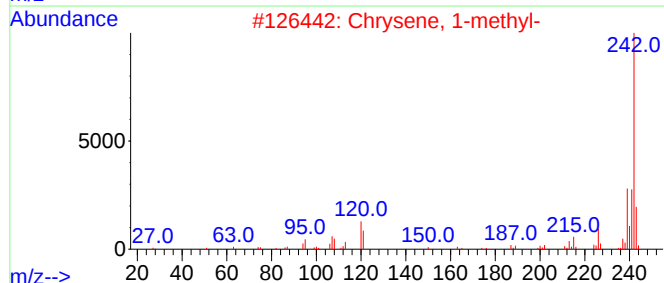
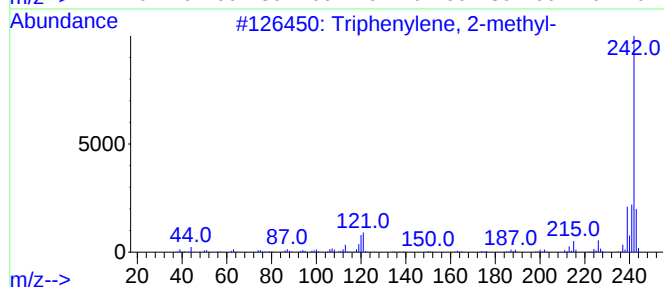
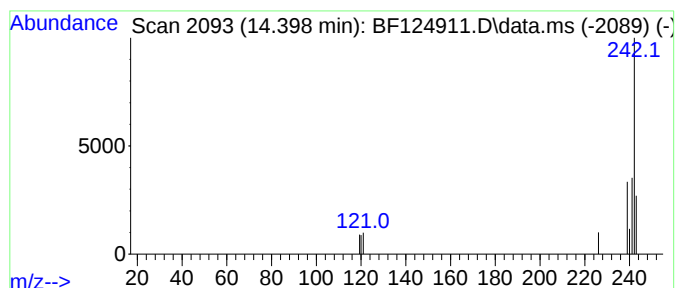
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 19 Triphenylene, 2-methyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.398	4.25 ng	9873	Chrysene-d12	14.009

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Triphenylene, 2-methyl-	242	C19H14	001705-84-6	80
2			Chrysene, 1-methyl-	242	C19H14	003351-28-8	80
3			Benz[a]anthracene, 7-methyl-	242	C19H14	002541-69-7	80
4			Chrysene, 5-methyl-	242	C19H14	003697-24-3	80
5			2-Methylchrysene	242	C19H14	003351-32-4	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF080221\
Data File : BF124911.D
Acq On : 2 Aug 2021 16:57
Operator : JU/CG
Sample : M3253-03
Misc :
ALS Vial : 15 Sample Multiplier: 1

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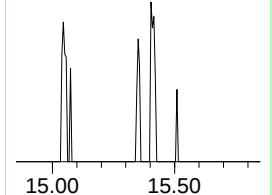
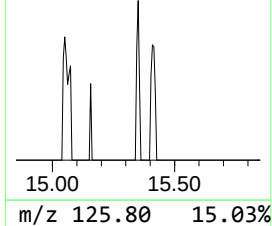
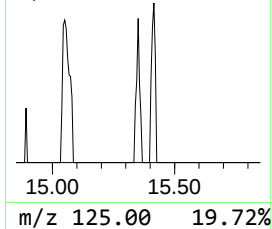
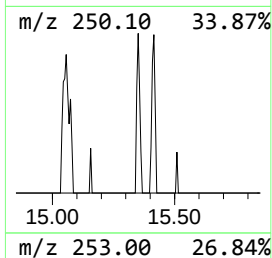
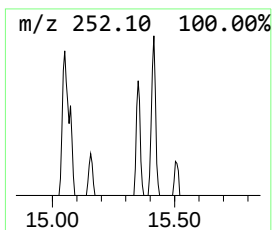
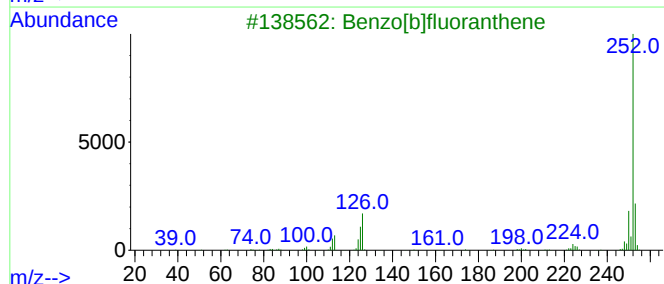
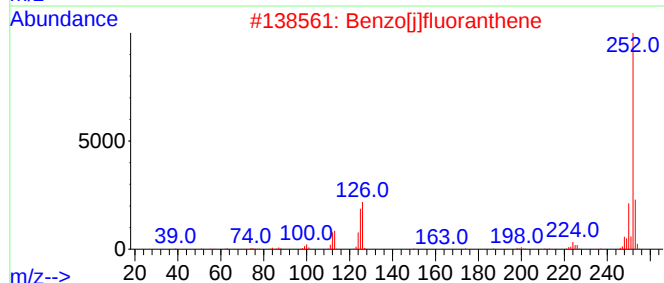
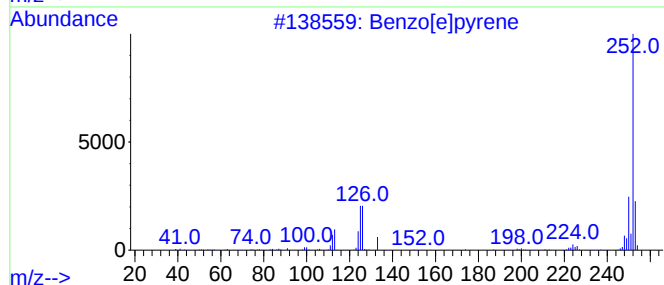
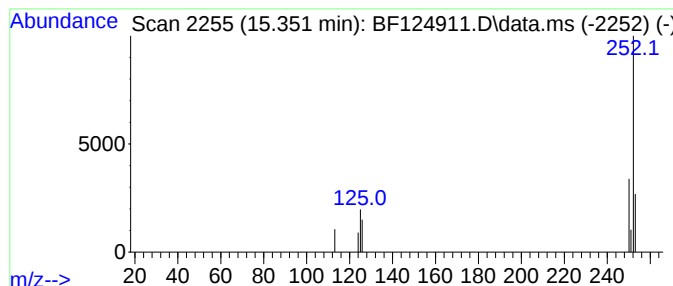
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 20 Benzo[e]pyrene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.351	5.72 ng	7438	Perylene-d12	15.480

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[e]pyrene	252	C20H12	000192-97-2	80
2			Benzo[j]fluoranthene	252	C20H12	000205-82-3	72
3			Benzo[b]fluoranthene	252	C20H12	000205-99-2	72
4			Perylene	252	C20H12	000198-55-0	72
5			Benzo[k]fluoranthene	252	C20H12	000207-08-9	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF080221\
 Data File : BF124911.D
 Acq On : 2 Aug 2021 16:57
 Operator : JU/CG
 Sample : M3253-03
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MH-2-COMP

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF072221.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
unknown2.116	2.116	11.9	ng	16717	1	6.851	27978	20.0
Naphthalene, 2-...	9.545	5.4	ng	11683	3	9.892	43257	20.0
9H-Fluorene, 2-...	10.980	4.2	ng	8454	4	11.374	40730	20.0
Phenanthrene, 1...	11.874	12.8	ng	25987	4	11.374	40730	20.0
Phenanthrene, 2...	11.904	13.8	ng	28151	4	11.374	40730	20.0
unknown11.986	11.986	14.6	ng	29668	4	11.374	40730	20.0
Phenanthrene, 4...	12.010	9.8	ng	20045	4	11.374	40730	20.0
Naphthalene, 2-...	12.168	3.0	ng	6067	4	11.374	40730	20.0
9,10-Dimethylan...	12.351	2.8	ng	5775	4	11.374	40730	20.0
di-p-Tolylacety...	12.427	8.7	ng	17766	4	11.374	40730	20.0
3-Amino-5-bromo...	12.463	7.0	ng	14230	4	11.374	40730	20.0
unknown12.510	12.510	4.4	ng	8875	4	11.374	40730	20.0
Fluoranthene, 2...	13.033	3.9	ng	9139	5	14.009	46467	20.0
11H-Benzo[a]flu...	13.145	8.8	ng	20428	5	14.009	46467	20.0
Pyrene, 1-methyl-	13.245	4.8	ng	11136	5	14.009	46467	20.0
Pyrene, 4-methyl-	13.339	4.3	ng	10095	5	14.009	46467	20.0
11H-Benzo[b]flu...	13.368	4.0	ng	9212	5	14.009	46467	20.0
5,6,8,9,10,11-H...	13.762	2.8	ng	6556	5	14.009	46467	20.0
Triphenylene, 2...	14.398	4.3	ng	9873	5	14.009	46467	20.0
Benzo[e]pyrene	15.351	5.7	ng	7438	6	15.480	25994	20.0