

Data Path : Z:\HPCHEM1\BNA F\DATA\BF081617\
 Data File : BF097755.D
 Acq On : 16 Aug 2017 22:54
 Operator : SJ/JU
 Sample : I4739-05 5X
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 S11-WC-5

Integration Parameters: rteint.p

Integrator: RTE

Smoothing : ON

Sampling : 1

Start Thrs: 0.2

Stop Thrs : 0

Filtering: 5

Min Area: 3 % of largest Peak

Max Peaks: 100

Peak Location: TOP

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

Peak separation: 5

Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF081517.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.963	486	489	494	rBV	70918	71243	4.22%	0.256%
2	5.375	555	559	563	rBV	1035032	895860	53.11%	3.224%
3	6.398	729	733	741	rVB	1189811	981967	58.22%	3.534%
4	6.539	753	757	761	rBV	1068068	955607	56.65%	3.439%
5	6.781	794	798	801	rBV	1304649	1072963	63.61%	3.862%
6	6.934	820	824	828	rBV	876031	694137	41.15%	2.498%
7	7.339	888	893	896	rBV	653348	571820	33.90%	2.058%
8	8.063	1011	1016	1019	rBV	1753390	1426326	84.56%	5.134%
9	8.775	1134	1137	1140	rVB	37773	26319	1.56%	0.095%
10	8.875	1150	1154	1157	rVB	48493	40392	2.39%	0.145%
11	9.139	1194	1199	1202	rBV	1093108	928566	55.05%	3.342%
12	9.475	1251	1256	1258	rBV	56016	51394	3.05%	0.185%
13	9.533	1263	1266	1270	rVB	126572	100603	5.96%	0.362%
14	9.675	1286	1290	1293	rBV	75910	61683	3.66%	0.222%
15	9.816	1307	1314	1317	rBV2	1648847	1419763	84.17%	5.110%
16	9.845	1317	1319	1323	rVB	201389	156373	9.27%	0.563%
17	10.016	1345	1348	1352	rVB	102801	86238	5.11%	0.310%
18	10.133	1362	1368	1370	rBV3	28796	32449	1.92%	0.117%
19	10.222	1378	1383	1385	rBV2	25143	29616	1.76%	0.107%
20	10.357	1403	1406	1413	rVB	131961	134044	7.95%	0.482%
21	10.427	1413	1418	1420	rBV	35636	44694	2.65%	0.161%
22	10.522	1431	1434	1438	rVB	54661	57385	3.40%	0.207%
23	10.598	1443	1447	1451	rVV	658503	550708	32.65%	1.982%
24	10.727	1464	1469	1473	rBV	64769	64674	3.83%	0.233%
25	10.910	1493	1500	1503	rBV4	44801	83171	4.93%	0.299%
26	10.933	1503	1504	1507	rVV3	33229	27422	1.63%	0.099%
27	11.039	1517	1522	1523	rVV2	53710	59578	3.53%	0.214%
28	11.057	1523	1525	1528	rVV2	52000	59093	3.50%	0.213%
29	11.104	1530	1533	1537	rVV	83382	86094	5.10%	0.310%
30	11.157	1537	1542	1544	rVV3	48205	76465	4.53%	0.275%
31	11.192	1546	1548	1554	rVB	91472	96113	5.70%	0.346%
32	11.298	1562	1566	1568	rBV	1303529	1226356	72.70%	4.414%
33	11.321	1568	1570	1574	rVV	1713887	1455736	86.30%	5.239%
34	11.374	1574	1579	1582	rVV	339960	310302	18.40%	1.117%

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Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF081517.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	11.404	1582	1584	1588	rVV2	38834	35946	2.13%	0.129%
36	11.527	1599	1605	1607	rBV2	119861	120696	7.16%	0.434%
37	11.598	1615	1617	1620	rBV2	65115	55027	3.26%	0.198%
38	11.627	1620	1622	1626	rVB2	38237	32610	1.93%	0.117%
39	11.798	1645	1651	1653	rBV	215189	191784	11.37%	0.690%
40	11.827	1653	1656	1660	rVB	294223	245300	14.54%	0.883%
41	11.874	1660	1664	1666	rBV	106835	101050	5.99%	0.364%
42	11.910	1666	1670	1672	rVV	308254	312335	18.52%	1.124%
43	11.933	1672	1674	1677	rVB	153634	115227	6.83%	0.415%
44	12.098	1699	1702	1703	rBV	144537	128984	7.65%	0.464%
45	12.116	1703	1705	1709	rVB	125723	97331	5.77%	0.350%
46	12.245	1724	1727	1730	rVB2	58612	45248	2.68%	0.163%
47	12.274	1730	1732	1734	rBV	41673	32777	1.94%	0.118%
48	12.298	1734	1736	1738	rVB2	56182	44451	2.64%	0.160%
49	12.345	1740	1744	1748	rBV	117294	141274	8.38%	0.508%
50	12.386	1748	1751	1753	rVV	63568	79489	4.71%	0.286%
51	12.427	1756	1758	1764	rVB3	94575	115207	6.83%	0.415%
52	12.510	1768	1772	1776	rVB	1774099	1521875	90.22%	5.477%
53	12.568	1776	1782	1785	rBV4	37078	65597	3.89%	0.236%
54	12.604	1785	1788	1790	rVB	38989	35984	2.13%	0.130%
55	12.639	1790	1794	1795	rBV3	43550	52228	3.10%	0.188%
56	12.657	1795	1797	1801	rVB	54231	57814	3.43%	0.208%
57	12.739	1805	1811	1814	rBV	1592519	1610441	95.48%	5.796%
58	12.786	1816	1819	1824	rVB2	73852	103496	6.14%	0.372%
59	12.857	1827	1831	1832	rBV2	83711	74133	4.39%	0.267%
60	12.880	1832	1835	1842	rVB	728589	711864	42.20%	2.562%
61	12.957	1842	1848	1851	rBV2	103332	174838	10.37%	0.629%
62	13.021	1855	1859	1860	rBV3	31788	38821	2.30%	0.140%
63	13.039	1860	1862	1863	rVV	71440	61360	3.64%	0.221%
64	13.063	1864	1866	1869	rVB	159358	135275	8.02%	0.487%
65	13.098	1869	1872	1874	rBV2	27081	33484	1.99%	0.121%
66	13.127	1874	1877	1880	rVB2	84937	79961	4.74%	0.288%
67	13.168	1880	1884	1887	rBV2	161911	165418	9.81%	0.595%
68	13.257	1897	1899	1902	rVB	97366	84398	5.00%	0.304%
69	13.286	1902	1904	1909	rBV	95061	109335	6.48%	0.394%
70	13.380	1914	1920	1927	rVB8	34925	82145	4.87%	0.296%
71	13.468	1927	1935	1937	rBV7	50869	105639	6.26%	0.380%

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72	13.568	1949	1952	1957	rVB	103499	113998	6.76%	0.410%
73	13.680	1966	1971	1974	rBV2	106046	164861	9.77%	0.593%
74	13.710	1974	1976	1978	rBV	96056	72689	4.31%	0.262%
75	13.733	1978	1980	1984	rVB2	64683	73033	4.33%	0.263%
76	13.774	1984	1987	1995	rVB2	125077	177747	10.54%	0.640%
77	13.933	2006	2014	2016	rBV2	1237195	1686760	100.00%	6.071%
78	13.957	2016	2018	2021	rVB	565265	449656	26.66%	1.618%
79	14.039	2029	2032	2035	rVB	71454	62524	3.71%	0.225%
80	14.115	2043	2045	2048	rVB2	55959	49942	2.96%	0.180%
81	14.151	2048	2051	2055	rVV2	44845	63780	3.78%	0.230%
82	14.251	2065	2068	2071	rBV4	32928	35029	2.08%	0.126%
83	14.280	2071	2073	2075	rBV	43189	40480	2.40%	0.146%
84	14.315	2076	2079	2083	rVV	122874	131117	7.77%	0.472%
85	14.351	2083	2085	2088	rVB	75253	62124	3.68%	0.224%
86	14.410	2093	2095	2098	rVV3	67328	74490	4.42%	0.268%
87	14.439	2098	2100	2102	rVV	59464	60892	3.61%	0.219%
88	14.462	2102	2104	2108	rVB3	38158	44953	2.67%	0.162%
89	14.615	2127	2130	2133	rBV2	35685	42875	2.54%	0.154%
90	14.698	2141	2144	2146	rBV4	33306	32784	1.94%	0.118%
91	14.780	2155	2158	2159	rBV2	30981	33087	1.96%	0.119%
92	14.951	2183	2187	2190	rBV	465691	694615	41.18%	2.500%
93	15.057	2201	2205	2208	rVB	104086	115981	6.88%	0.417%
94	15.245	2232	2237	2242	rVB	294329	356808	21.15%	1.284%
95	15.304	2242	2247	2253	rVB	410329	474897	28.15%	1.709%
96	15.362	2253	2257	2261	rBV	622838	823400	48.82%	2.964%
97	15.392	2261	2262	2270	rVB2	130927	153908	9.12%	0.554%
98	16.327	2419	2421	2428	rVB6	26293	46481	2.76%	0.167%
99	16.756	2489	2494	2502	rVB2	149244	287055	17.02%	1.033%
100	17.174	2560	2565	2576	rVB	110426	224442	13.31%	0.808%

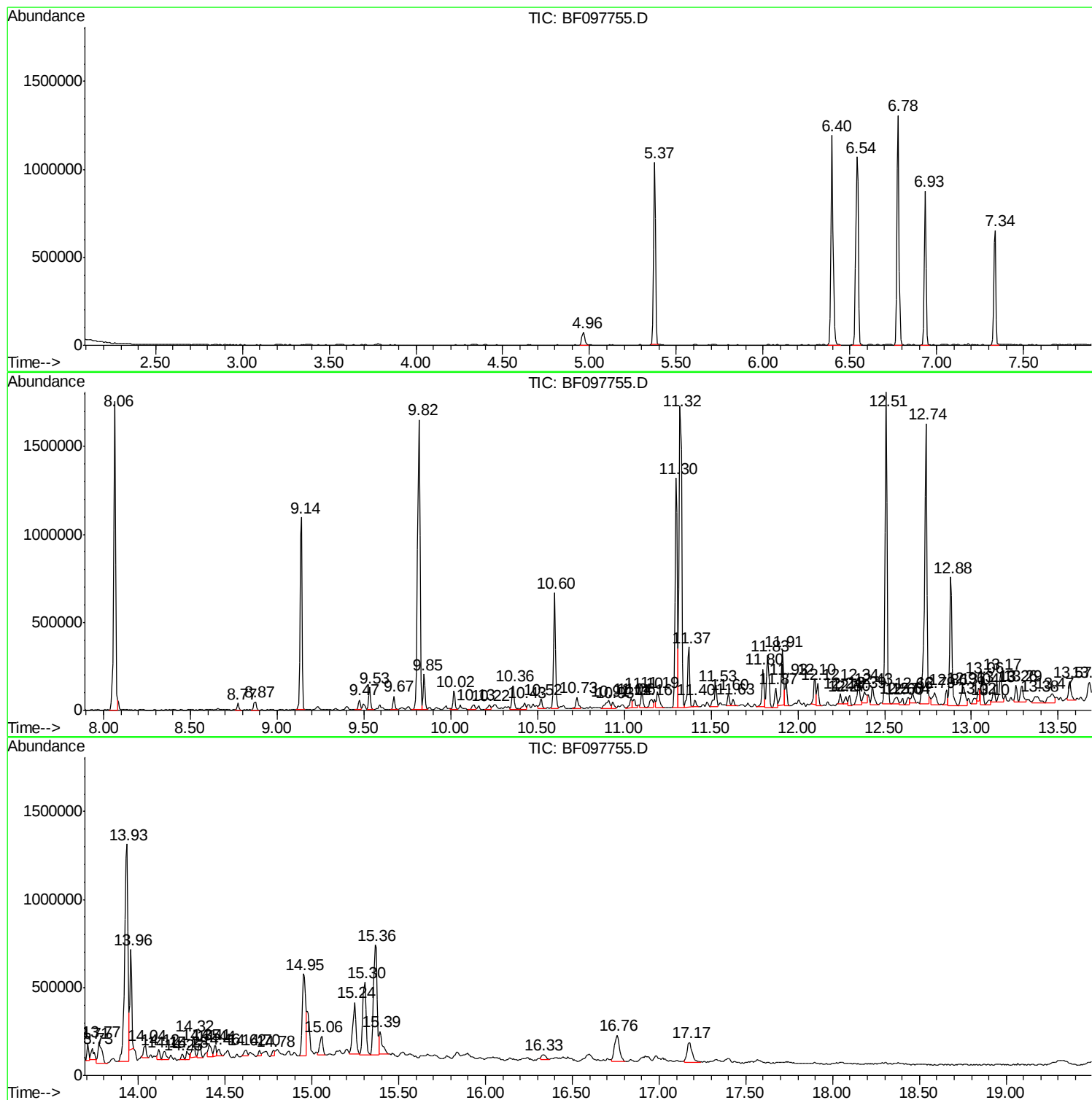
Sum of corrected areas: 27784404

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF081517.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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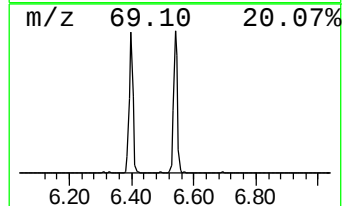
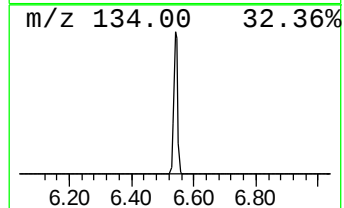
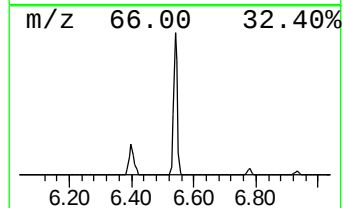
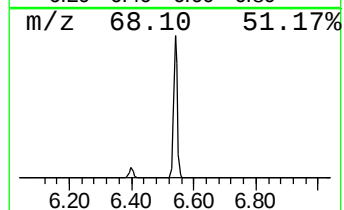
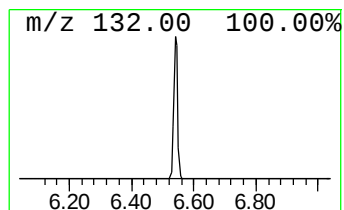
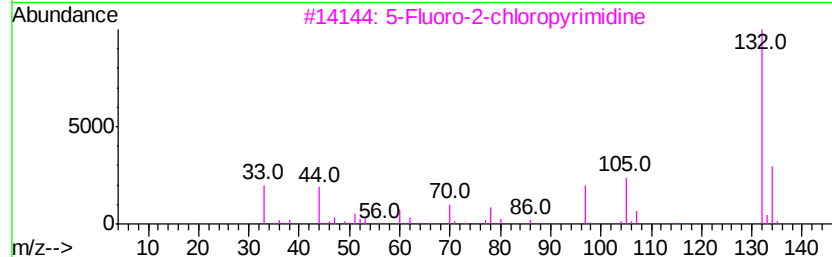
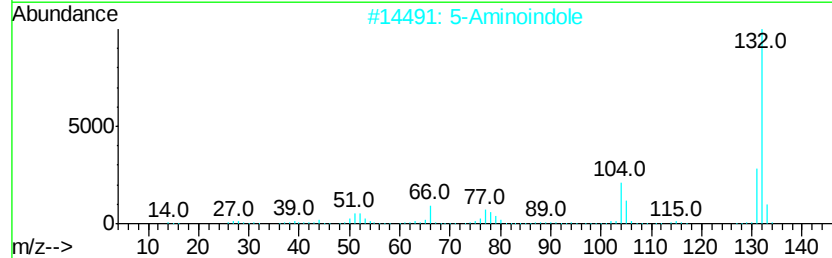
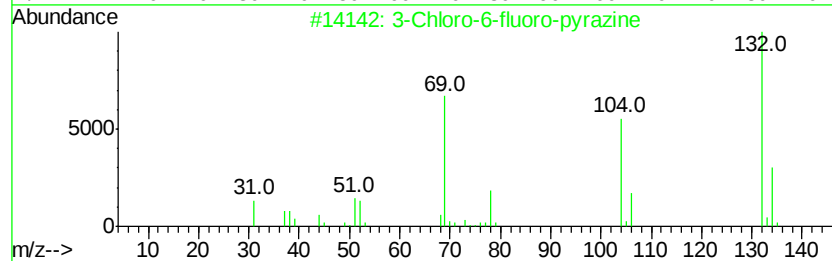
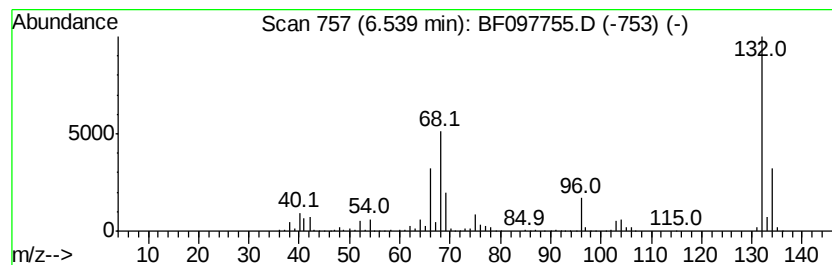
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF081517.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 unknown6.54 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.54	17.81 ng	955607	1,4-Dichlorobenzene-d4	6.78

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2			5-Aminoindole	132	C8H8N2	005192-03-0	27
3			5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
4			2-Cyclohexen-1-one	96	C6H8O	000930-68-7	10
5			Cyclopropanamine, 1-phenyl-	133	C9H11N	1000351-26-2	10



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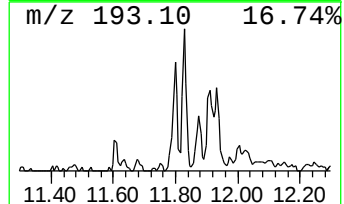
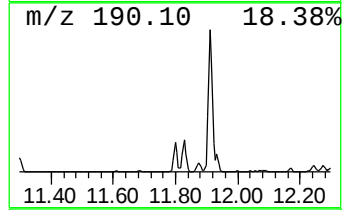
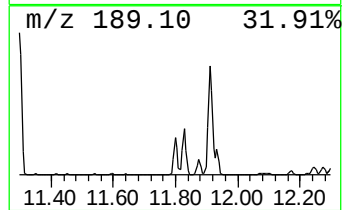
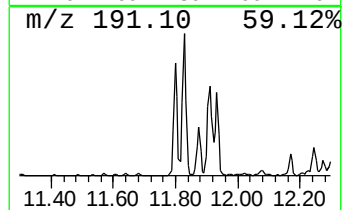
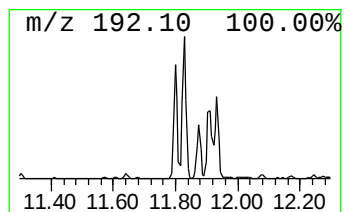
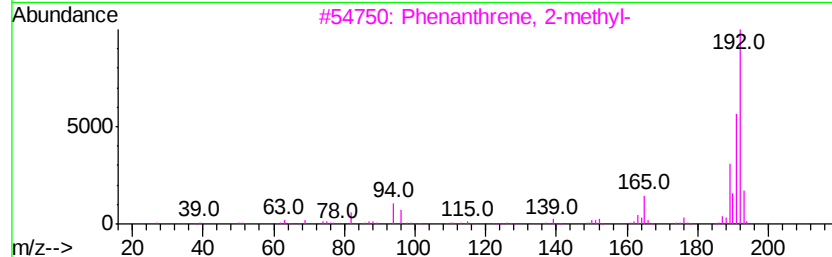
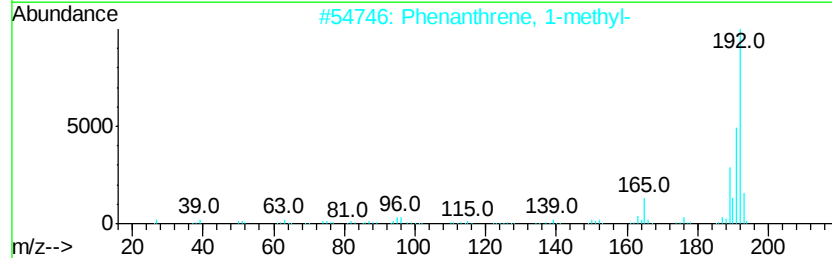
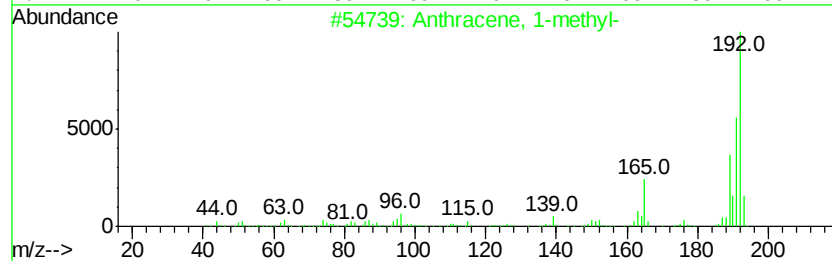
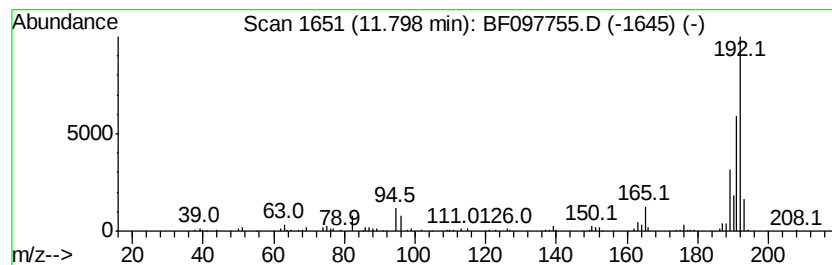
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Anthracene, 1-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.80	3.13 ng	191784	Phenanthrene-d10	11.30

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Anthracene, 1-methyl-	192	C15H12	000610-48-0	94
2		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	93
3		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	93
4		Anthracene, 2-methyl-	192	C15H12	000613-12-7	90
5		Anthracene, 9-methyl-	192	C15H12	000779-02-2	87



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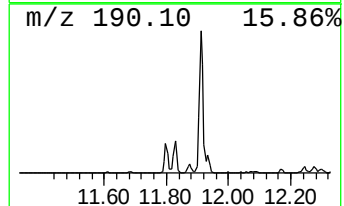
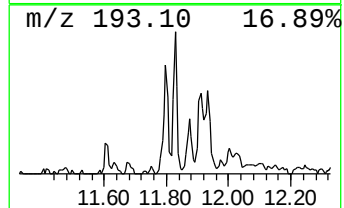
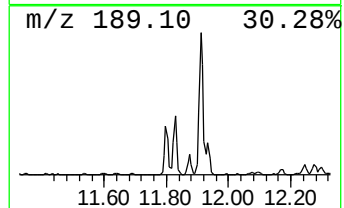
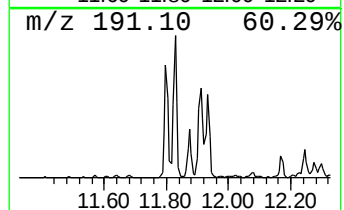
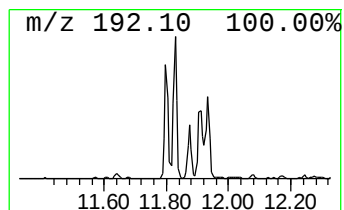
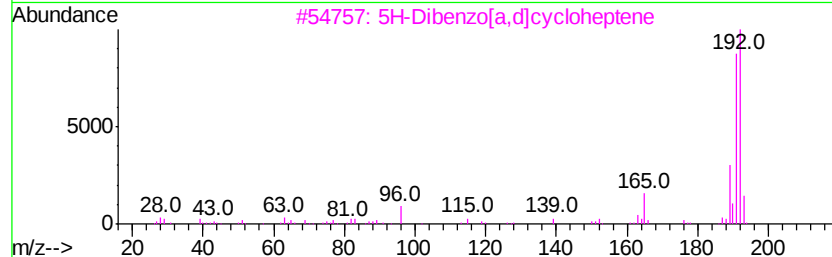
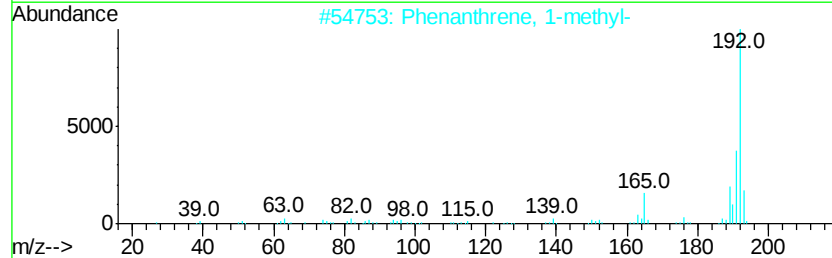
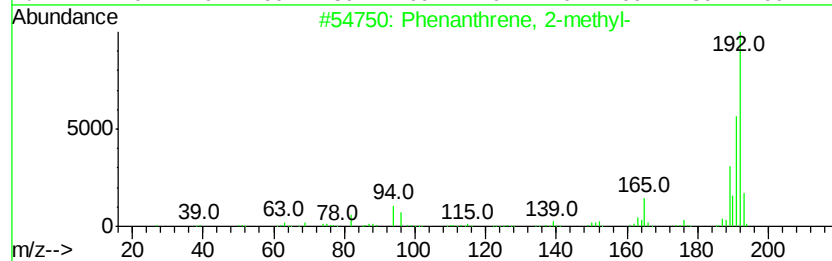
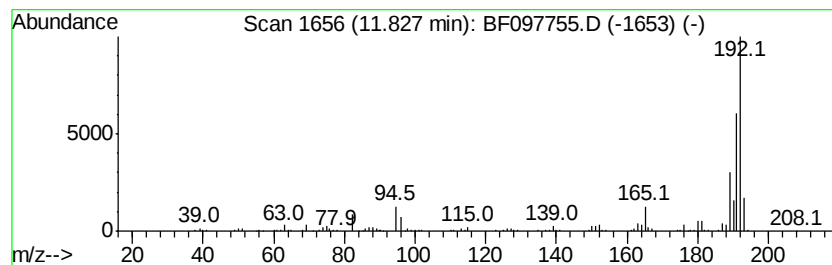
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TIC Integration Parameters: LSCINT.P

Peak Number 4 Phenanthrene, 2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.83	4.00 ng	245300	Phenanthrene-d10	11.30	
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	95
3	5H-Dibenzo[a,d]cycloheptene	192	C15H12	000256-81-5	93
4	Phenanthrene, 3-methyl-	192	C15H12	000832-71-3	93
5	Anthracene, 2-methyl-	192	C15H12	000613-12-7	93



Data Path : Z:\HPCHEM1\BNA F\DATA\BF081617\
Data File : BF097755.D
Acq On : 16 Aug 2017 22:54
Operator : SJ/JU
Sample : I4739-05 5X
Misc :
ALS Vial : 23 Sample Multiplier: 1

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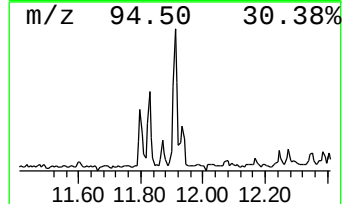
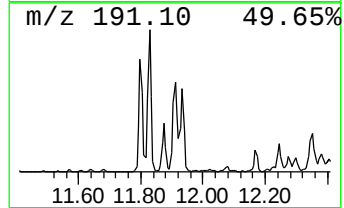
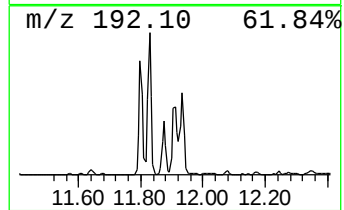
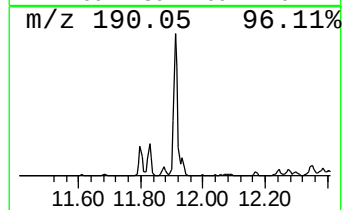
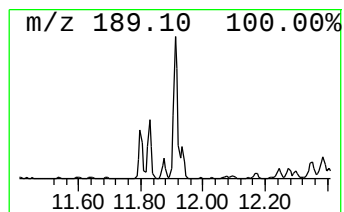
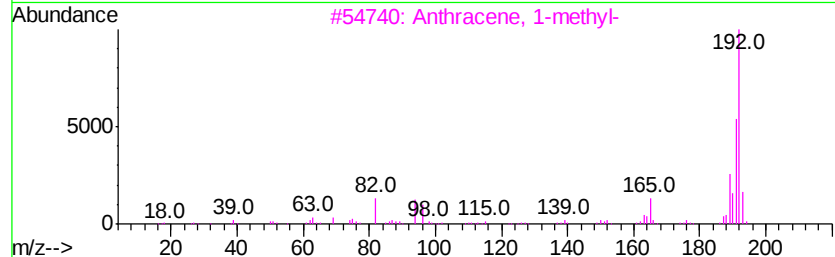
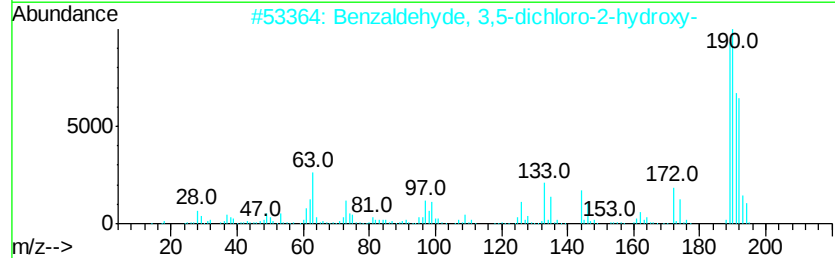
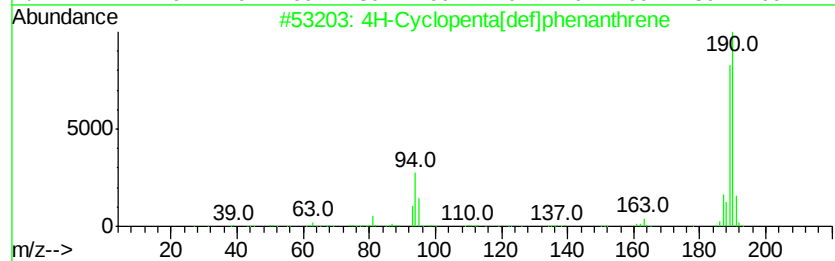
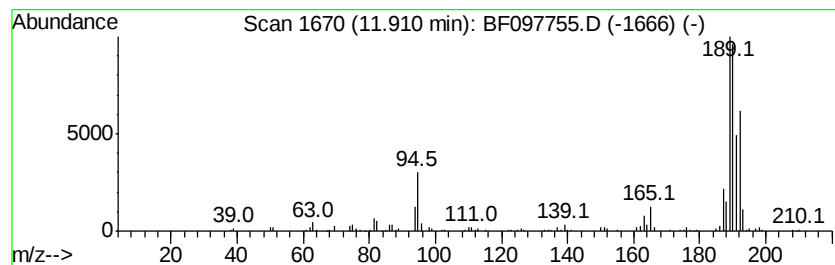
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 4H-Cyclopenta[def]phenanthrene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.91	5.09 ng	312335	Phenanthrene-d10	11.30

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	64
2		Benzaldehyde, 3,5-dichloro-2-hyd...	190	C7H4Cl2O2	000090-60-8	64
3		Anthracene, 1-methyl-	192	C15H12	000610-48-0	18
4		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	18
5		1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	18



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF081617\
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Misc :
ALS Vial : 23 Sample Multiplier: 1

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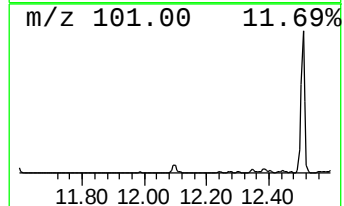
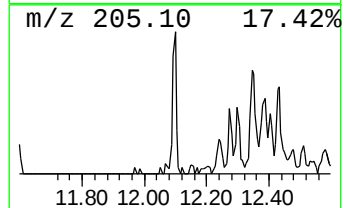
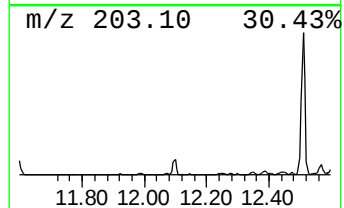
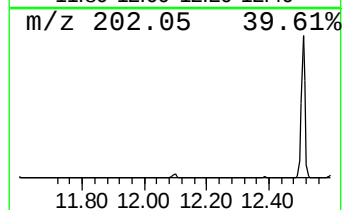
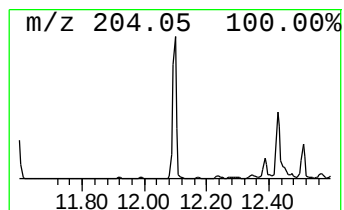
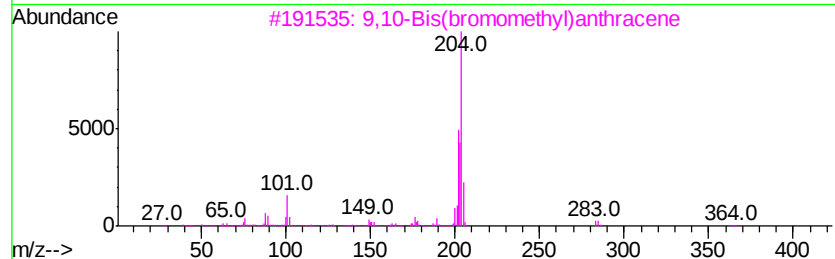
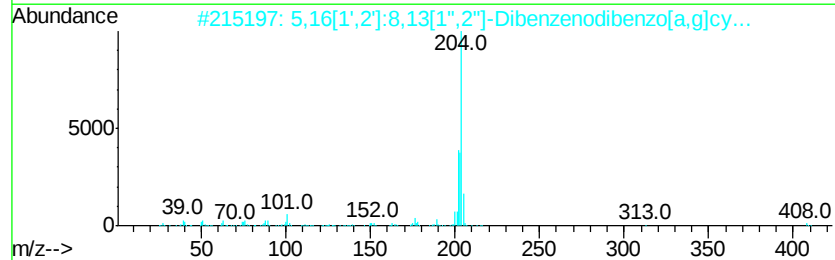
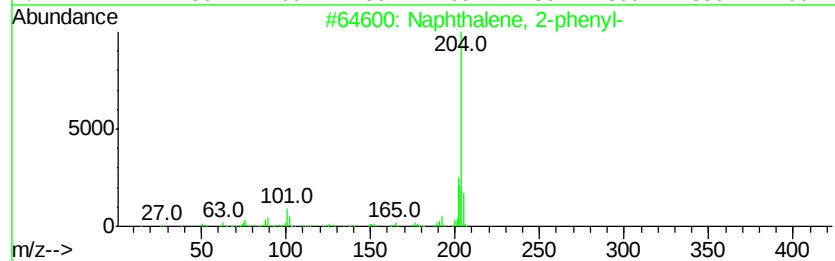
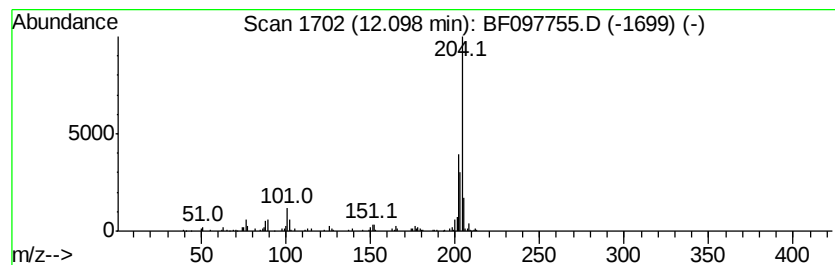
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Naphthalene, 2-phenyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.10	2.10 ng	128984	Phenanthrene-d10	11.30

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	92
2		5,16[1',2']:8,13[1'',2'']-Dibenz...	408	C32H24	005672-97-9	86
3		9,10-Bis(bromomethyl)anthracene	362	C16H12Br2	034373-96-1	70
4		Tricyclo[8.2.2.2(4,7)]hexadeca-2...	204	C16H12	006572-60-7	70
5		Naphthalene, 1-phenyl-	204	C16H12	000605-02-7	55



Data Path : Z:\HPCHEM1\BNA F\DATA\BF081617\
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ALS Vial : 23 Sample Multiplier: 1

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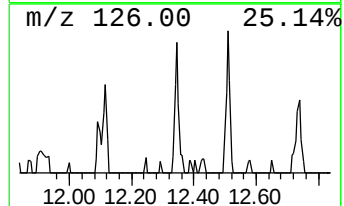
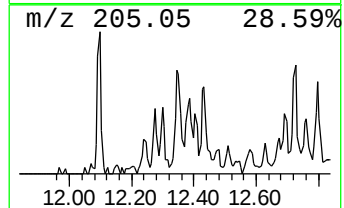
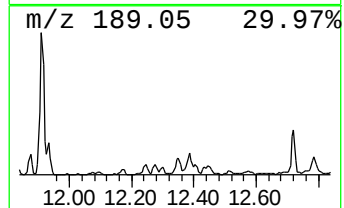
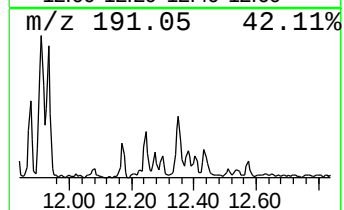
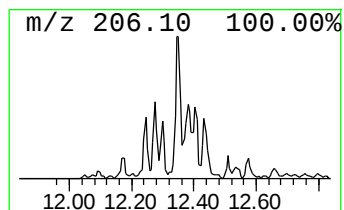
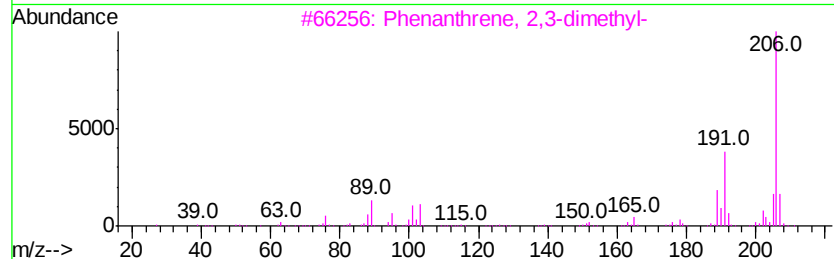
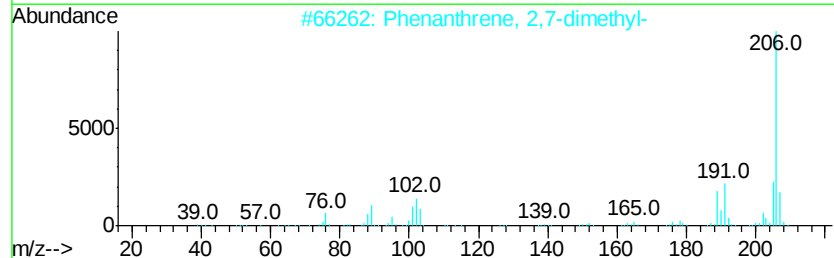
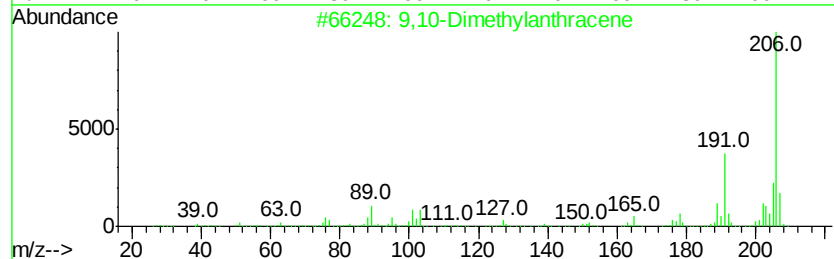
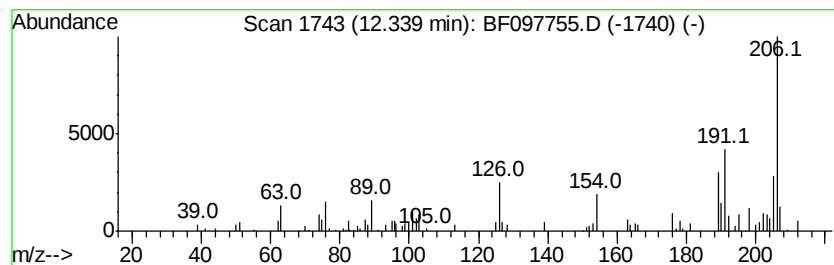
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 9,10-Dimethylantracene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD		R.T.
12.34	2.30 ng	141274	Phenanthrene-d10		11.30

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	9,10-Dimethylantracene			206	C16H14	000781-43-1	93
2	Phenanthrene, 2,7-dimethyl-			206	C16H14	001576-69-8	90
3	Phenanthrene, 2,3-dimethyl-			206	C16H14	003674-65-5	90
4	Phenanthrene, 3,6-dimethyl-			206	C16H14	001576-67-6	87
5	Phenanthrene, 2,5-dimethyl-			206	C16H14	003674-66-6	83



Data Path : Z:\HPCHEM1\BNA F\DATA\BF081617\
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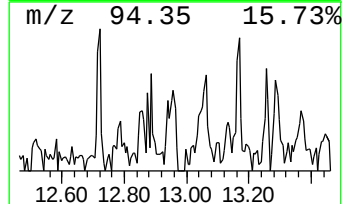
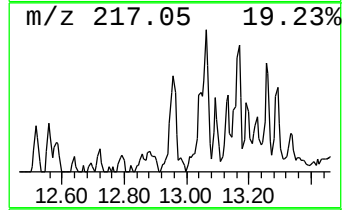
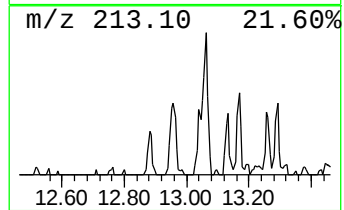
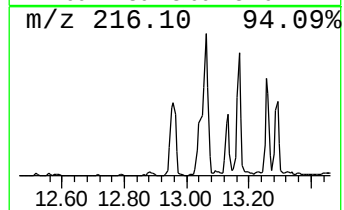
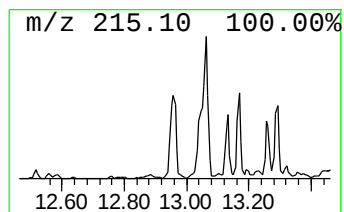
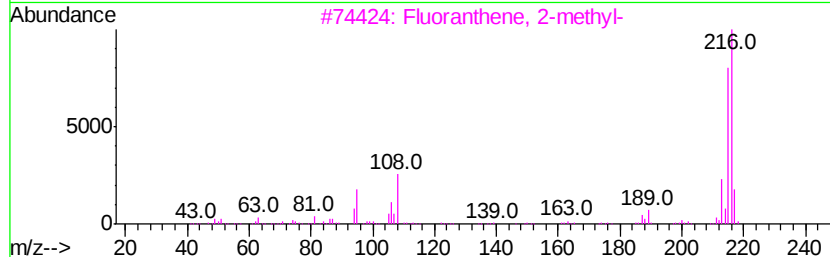
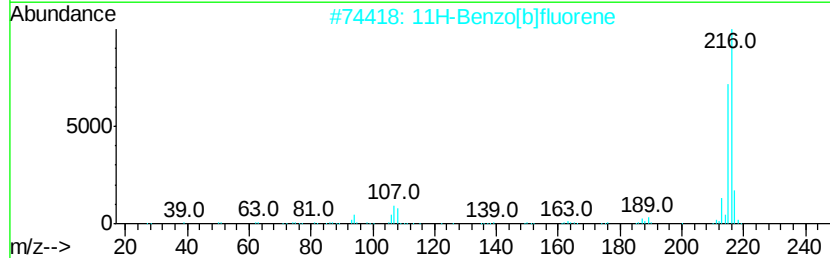
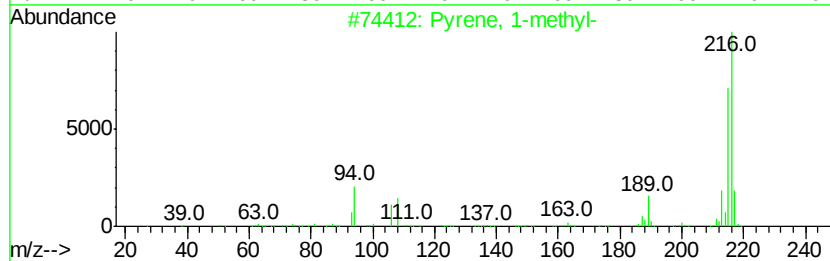
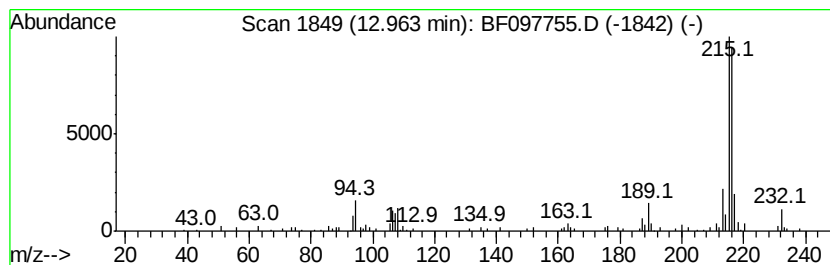
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Pyrene, 1-methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.96	2.07 ng	174838	Chrysene-d12	13.93
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Pyrene, 1-methyl-	216 C17H12	002381-21-7 98
2		11H-Benzo[b]fluorene	216 C17H12	000243-17-4 87
3		Fluoranthene, 2-methyl-	216 C17H12	033543-31-6 76
4		7H-Benzo[c]fluorene	216 C17H12	000205-12-9 64
5		Pyrene, 2-methyl-	216 C17H12	003442-78-2 62



Data Path : Z:\HPCHEM1\BNA F\DATA\BF081617\
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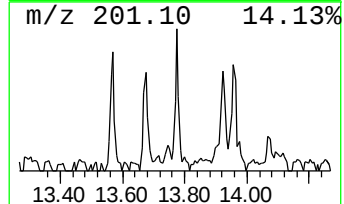
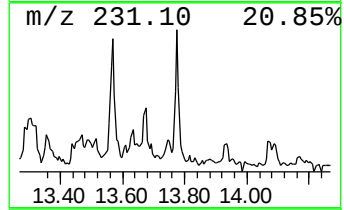
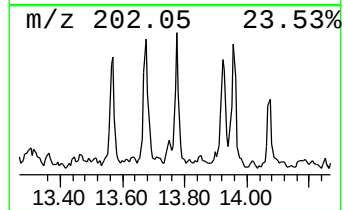
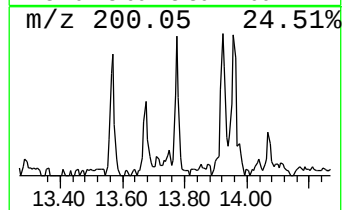
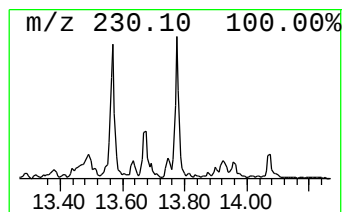
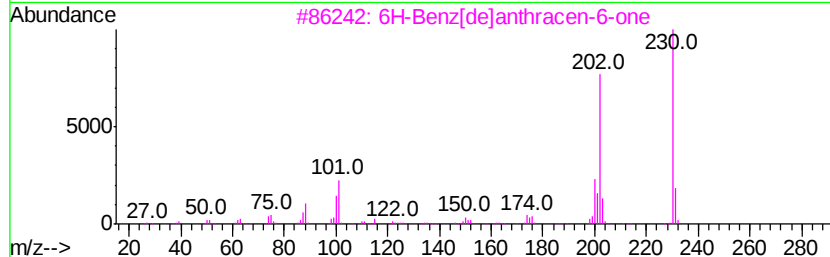
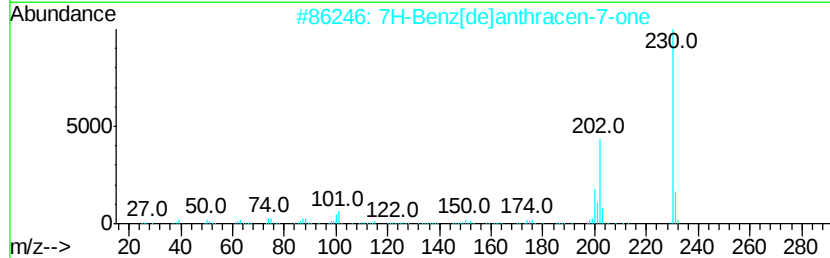
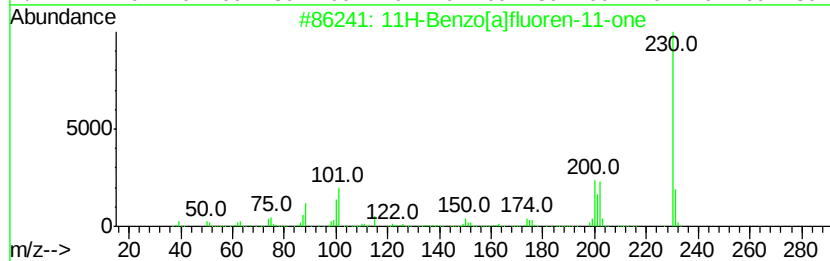
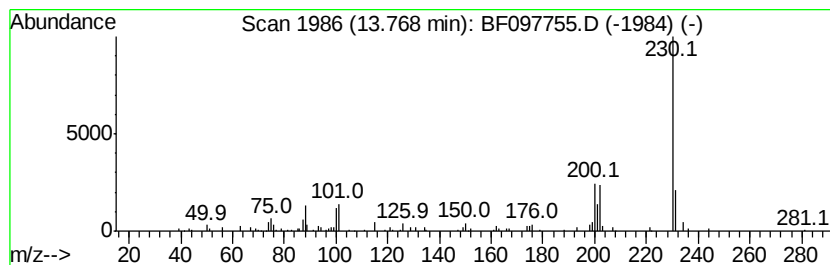
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 11H-Benzo[a]fluoren-11-one Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.77	2.11 ng	177747	Chrysene-d12	13.93

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	11H-Benzo[a]fluoren-11-one	230	C17H10O	000479-79-8	94
2		7H-Benz[de]anthracen-7-one	230	C17H10O	000082-05-3	91
3		6H-Benz[de]anthracen-6-one	230	C17H10O	080252-14-8	64
4		o-Terphenyl	230	C18H14	000084-15-1	59
5		5,6-Dihydrochrysene	230	C18H14	002091-92-1	52



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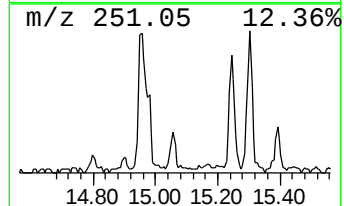
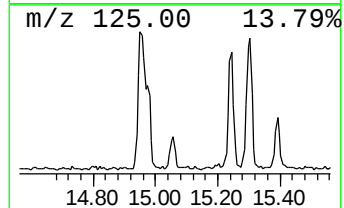
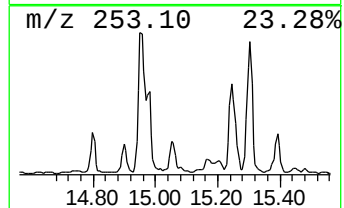
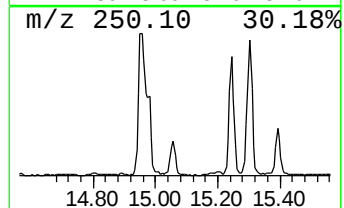
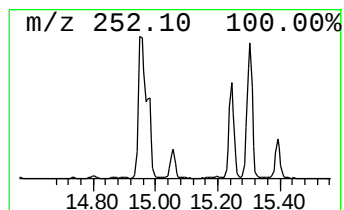
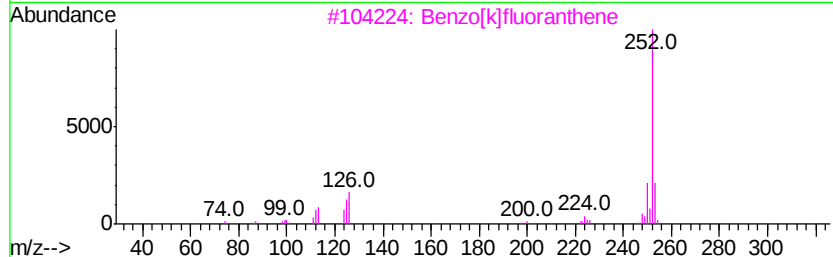
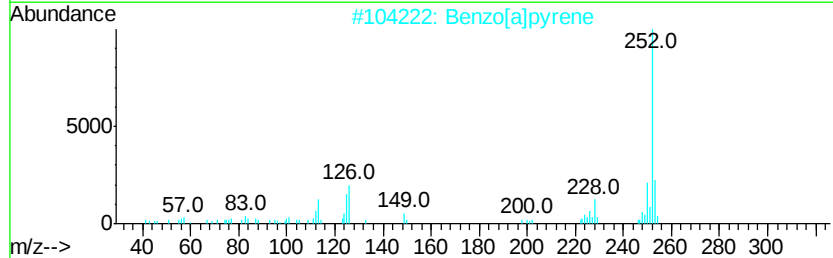
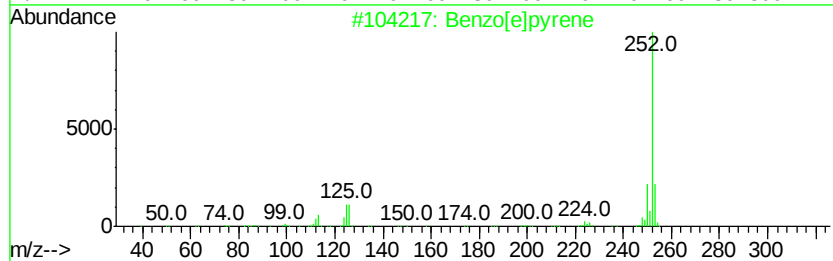
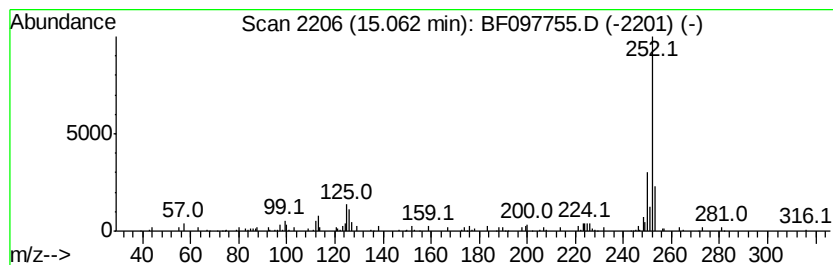
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Benzo[e]pyrene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.06	2.82 ng	115981	Perylene-d12	15.37

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



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Instrument :
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF081517.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
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Phenanthrene, 2-m...	11.83	4.0	ng	245300	4	11.30	1226360	20.0
4H-Cyclopenta[def...	11.91	5.1	ng	312335	4	11.30	1226360	20.0
Naphthalene, 2-ph...	12.10	2.1	ng	128984	4	11.30	1226360	20.0
9,10-Dimethylanthe...	12.34	2.3	ng	141274	4	11.30	1226360	20.0
Pyrene, 1-methyl-	12.96	2.1	ng	174838	5	13.93	1686760	20.0
11H-Benzo[a]fluor...	13.77	2.1	ng	177747	5	13.93	1686760	20.0
Benzo[e]pyrene	15.06	2.8	ng	115981	6	15.37	823400	20.0