

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF022020\
 Data File : BF118974.D
 Acq On : 20 Feb 2020 22:42
 Operator : CG/JU
 Sample : L1546-03
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 INDIAN-ISLAND-(2)

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:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF022020.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	5.387	557	561	576	rBV	3832271	4071499	81.72%	4.012%
2	6.404	729	734	737	rBV	3717448	3932150	78.93%	3.875%
3	6.757	791	794	798	rBV	1267024	1102762	22.13%	1.087%
4	6.916	817	821	825	rBV	3974680	3581594	71.89%	3.529%
5	7.322	882	890	893	rBV	2658044	2669022	53.57%	2.630%
6	8.039	1007	1012	1015	rBV	1635381	1444295	28.99%	1.423%
7	9.116	1189	1195	1198	rBV	5649239	4982115	100.00%	4.909%
8	9.380	1234	1240	1243	rBV	367185	490584	9.85%	0.483%
9	9.451	1248	1252	1254	rBV	826241	747692	15.01%	0.737%
10	9.475	1254	1256	1259	rVB	482789	369937	7.43%	0.365%
11	9.510	1259	1262	1268	rVB3	368316	454682	9.13%	0.448%
12	9.569	1268	1272	1274	rBV	390047	404797	8.13%	0.399%
13	9.651	1281	1286	1291	rBV	1064484	898740	18.04%	0.886%
14	9.792	1303	1310	1313	rVV	1674989	1668233	33.48%	1.644%
15	9.898	1324	1328	1332	rBV2	265460	351113	7.05%	0.346%
16	9.992	1340	1344	1347	rVV	509866	554882	11.14%	0.547%
17	10.033	1347	1351	1354	rVV	620080	519067	10.42%	0.511%
18	10.104	1359	1363	1366	rBV2	545272	624624	12.54%	0.615%
19	10.133	1366	1368	1371	rVV	346623	314752	6.32%	0.310%
20	10.198	1374	1379	1380	rBV2	341216	452425	9.08%	0.446%
21	10.286	1389	1394	1396	rVV4	278868	336709	6.76%	0.332%
22	10.333	1399	1402	1409	rVV3	539686	769564	15.45%	0.758%
23	10.398	1409	1413	1415	rVV2	244003	355285	7.13%	0.350%
24	10.451	1419	1422	1426	rVV4	239874	346725	6.96%	0.342%
25	10.492	1426	1429	1435	rVV4	280417	472585	9.49%	0.466%
26	10.580	1438	1444	1447	rVV	3046059	3112038	62.46%	3.066%
27	10.704	1461	1465	1470	rVB2	366580	380076	7.63%	0.375%
28	10.880	1491	1495	1498	rVV3	463309	709922	14.25%	0.700%
29	10.910	1498	1500	1503	rVV	505519	477519	9.58%	0.471%
30	10.969	1507	1510	1512	rVV	315741	335458	6.73%	0.331%
31	11.016	1515	1518	1519	rVV	239662	281904	5.66%	0.278%
32	11.033	1519	1521	1523	rVV2	281434	273299	5.49%	0.269%
33	11.080	1526	1529	1533	rVV2	309229	381269	7.65%	0.376%
34	11.169	1541	1544	1549	rVB	403116	450175	9.04%	0.444%

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Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF022020.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	11.274	1558	1562	1563	rBV	1319434	1249799	25.09%	1.231%
36	11.298	1563	1566	1569	rVV	3154800	2689088	53.97%	2.650%
37	11.345	1571	1574	1577	rVB	778445	662460	13.30%	0.653%
38	11.380	1577	1580	1583	rBV2	363265	431018	8.65%	0.425%
39	11.604	1614	1618	1622	rBV2	764451	733812	14.73%	0.723%
40	11.686	1627	1632	1636	rVB2	438495	584546	11.73%	0.576%
41	11.727	1636	1639	1642	rBV2	248248	270503	5.43%	0.267%
42	11.774	1643	1647	1649	rBV	1753030	1591588	31.95%	1.568%
43	11.804	1649	1652	1657	rVB	2398791	2096748	42.09%	2.066%
44	11.851	1657	1660	1662	rBV	637300	559510	11.23%	0.551%
45	11.886	1662	1666	1668	rVV	2111351	2223026	44.62%	2.190%
46	11.910	1668	1670	1675	rVB	1749235	1396220	28.02%	1.376%
47	12.010	1684	1687	1689	rVB	373236	314444	6.31%	0.310%
48	12.069	1689	1697	1699	rBV2	812519	977942	19.63%	0.964%
49	12.098	1699	1702	1707	rVB2	593219	730497	14.66%	0.720%
50	12.198	1717	1719	1721	rBV	337324	351605	7.06%	0.346%
51	12.216	1721	1722	1725	rVV2	383539	336271	6.75%	0.331%
52	12.251	1725	1728	1730	rVV	530753	488569	9.81%	0.481%
53	12.274	1730	1732	1734	rVB2	429170	329423	6.61%	0.325%
54	12.327	1734	1741	1743	rBV	1669556	1813851	36.41%	1.787%
55	12.363	1743	1747	1749	rVV	904414	1187040	23.83%	1.170%
56	12.410	1752	1755	1760	rBV2	718555	949878	19.07%	0.936%
57	12.486	1763	1768	1771	rBV	2948102	2580587	51.80%	2.543%
58	12.527	1772	1775	1777	rBV	335947	319515	6.41%	0.315%
59	12.580	1782	1784	1787	rVB2	347759	316733	6.36%	0.312%
60	12.716	1801	1807	1810	rBV	3614243	3955762	79.40%	3.898%
61	12.774	1813	1817	1820	rVB3	454098	552253	11.08%	0.544%
62	12.857	1828	1831	1838	rVB	4140363	4099844	82.29%	4.040%
63	12.933	1840	1844	1851	rBV3	813600	1271435	25.52%	1.253%
64	12.986	1851	1853	1856	rBV	352673	338164	6.79%	0.333%
65	13.021	1856	1859	1860	rBV2	616504	662911	13.31%	0.653%
66	13.045	1860	1863	1865	rVB	1009837	1052566	21.13%	1.037%
67	13.086	1865	1870	1871	rBV2	248674	273682	5.49%	0.270%
68	13.110	1871	1874	1877	rVB2	508906	439591	8.82%	0.433%
69	13.145	1877	1880	1884	rBV	1073324	954556	19.16%	0.941%
70	13.239	1892	1896	1898	rVB	903698	742584	14.90%	0.732%
71	13.268	1898	1901	1904	rVB	1072501	917390	18.41%	0.904%

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72	13.551	1946	1949	1953	rVB	589354	678729	13.62%	0.669%
73	13.657	1962	1967	1970	rBV2	706406	958768	19.24%	0.945%
74	13.686	1970	1972	1974	rVB	415817	312925	6.28%	0.308%
75	13.757	1980	1984	1988	rVB3	729472	919751	18.46%	0.906%
76	13.904	2004	2009	2012	rBV2	2262350	3202465	64.28%	3.156%
77	13.939	2012	2015	2022	rVB	1807342	1953935	39.22%	1.925%
78	13.998	2022	2025	2027	rBV	285140	281178	5.64%	0.277%
79	14.168	2051	2054	2057	rVV3	348189	431977	8.67%	0.426%
80	14.227	2061	2064	2067	rVV3	261187	291656	5.85%	0.287%
81	14.262	2067	2070	2071	rVV	278351	282408	5.67%	0.278%
82	14.298	2071	2076	2079	rVV	1124117	1418030	28.46%	1.397%
83	14.327	2079	2081	2084	rVV	722949	718545	14.42%	0.708%
84	14.374	2085	2089	2090	rVV3	420337	609168	12.23%	0.600%
85	14.392	2090	2092	2094	rVV2	545020	480164	9.64%	0.473%
86	14.421	2094	2097	2099	rVV	803415	774233	15.54%	0.763%
87	14.439	2099	2100	2104	rVV2	391194	341280	6.85%	0.336%
88	14.486	2104	2108	2112	rVV2	309279	421420	8.46%	0.415%
89	14.674	2138	2140	2142	rBV	298998	290016	5.82%	0.286%
90	14.710	2144	2146	2150	rVB	289044	300946	6.04%	0.297%
91	14.780	2155	2158	2161	rVV2	290762	358559	7.20%	0.353%
92	14.833	2165	2167	2174	rVB4	209243	297917	5.98%	0.294%
93	14.933	2179	2184	2190	rBV	1079886	1832722	36.79%	1.806%
94	15.033	2198	2201	2206	rVB	290733	317935	6.38%	0.313%
95	15.215	2228	2232	2238	rVB	768673	976995	19.61%	0.963%
96	15.280	2238	2243	2247	rBV	1147145	1456315	29.23%	1.435%
97	15.339	2249	2253	2256	rBV	884931	1097361	22.03%	1.081%
98	15.504	2276	2281	2284	rBV3	164917	326084	6.55%	0.321%
99	16.739	2485	2491	2498	rVB	383833	736257	14.78%	0.725%
100	17.156	2557	2562	2571	rVB	326306	657127	13.19%	0.647%

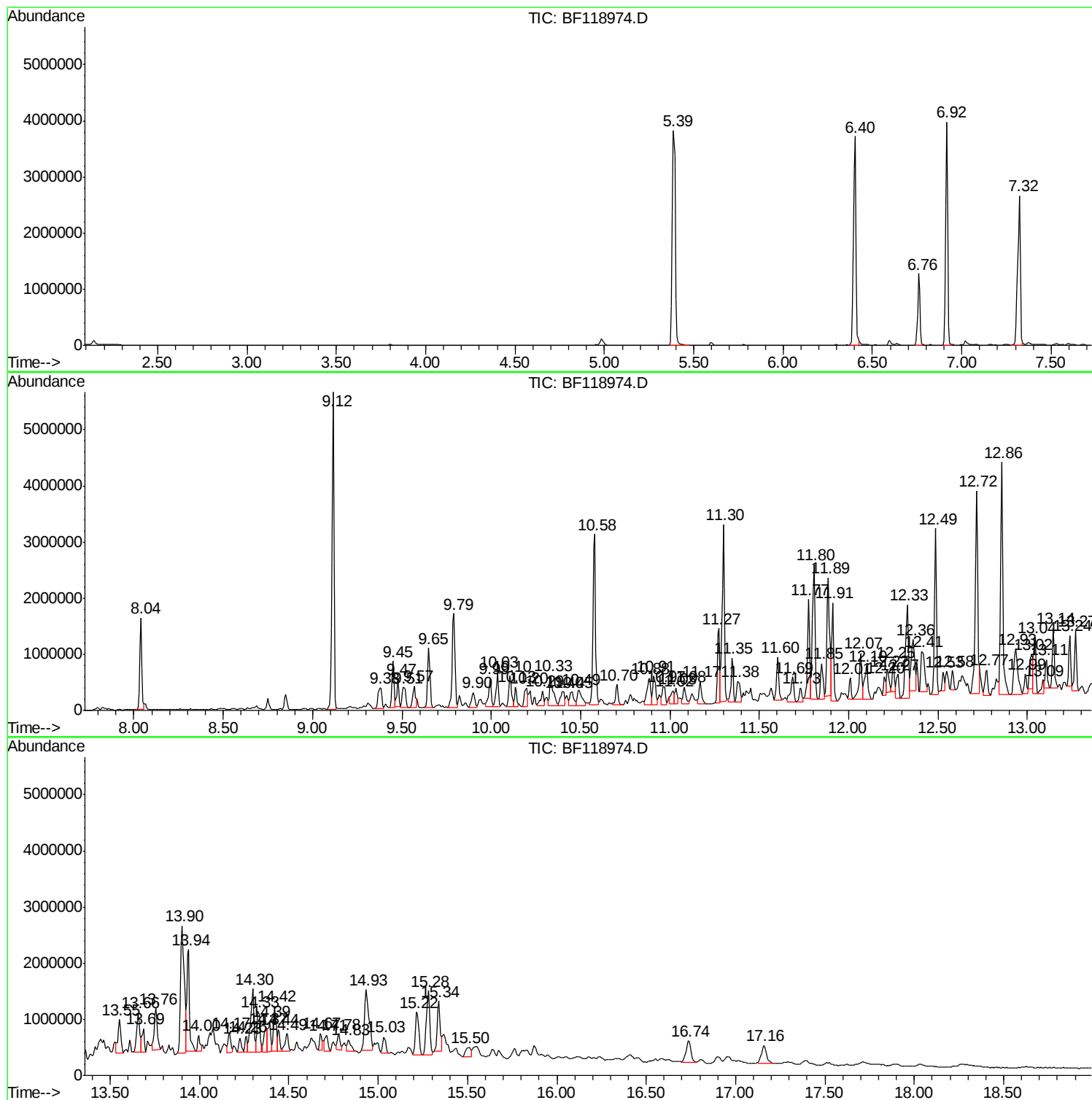
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INDIAN-ISLAND-(2)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022020.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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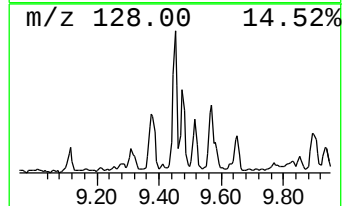
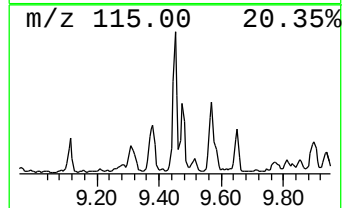
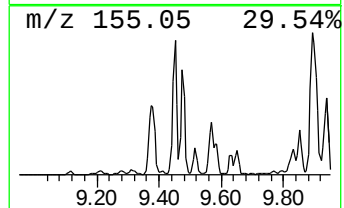
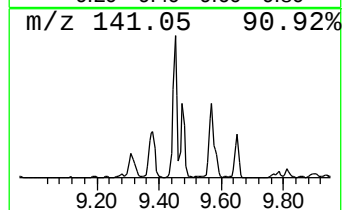
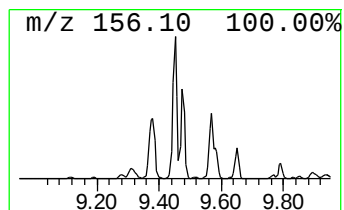
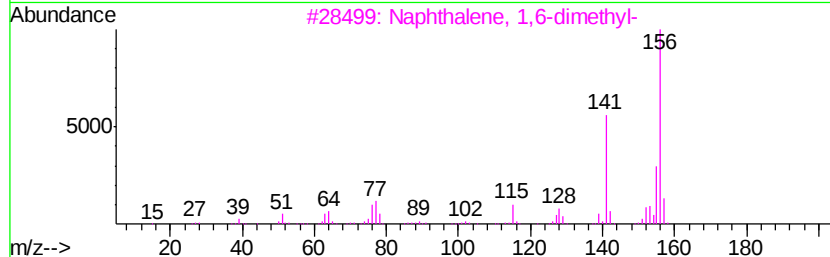
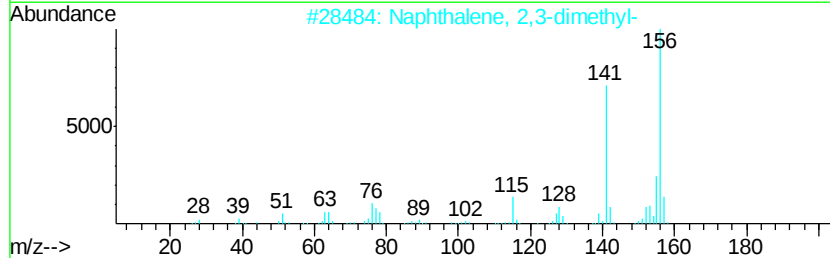
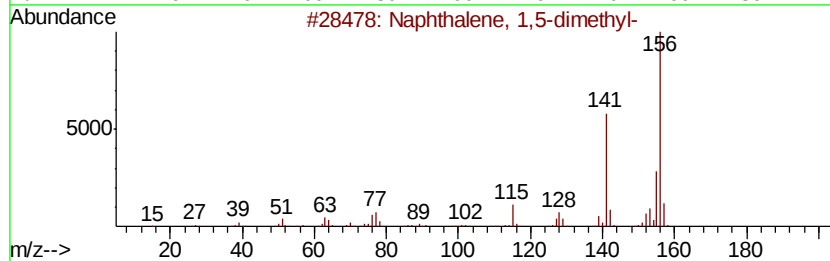
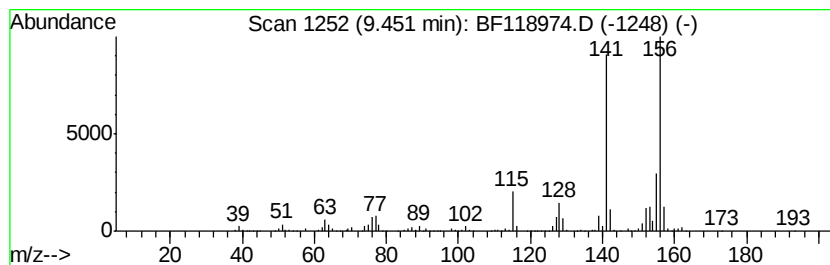
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 Naphthalene, 1,5-dimethyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.45	8.96 ng	747692	Acenaphthene-d10	9.79

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	97
2			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
3			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	97
4			Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	97
5			Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	97



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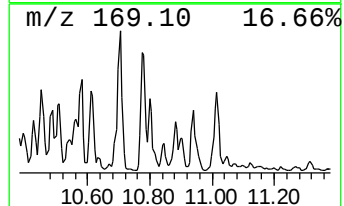
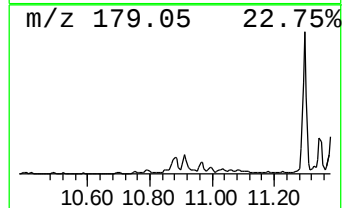
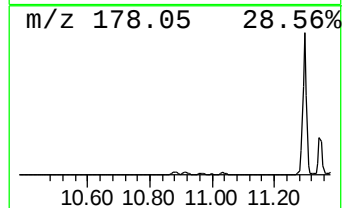
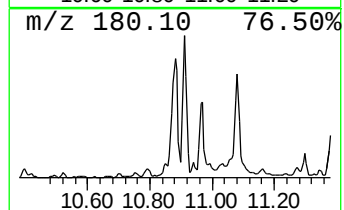
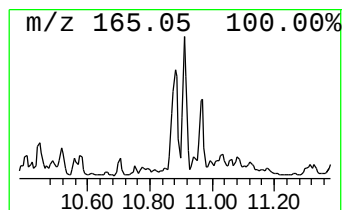
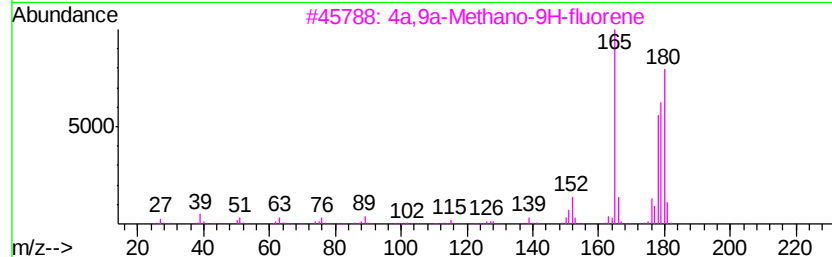
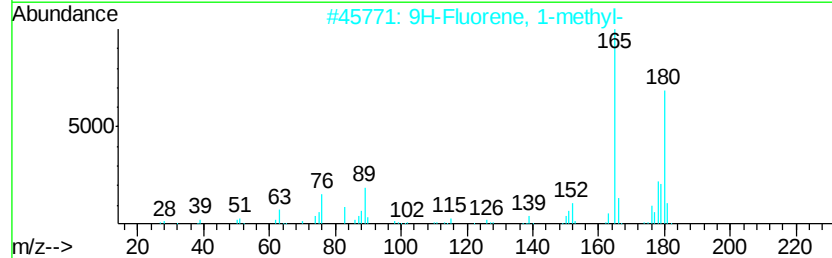
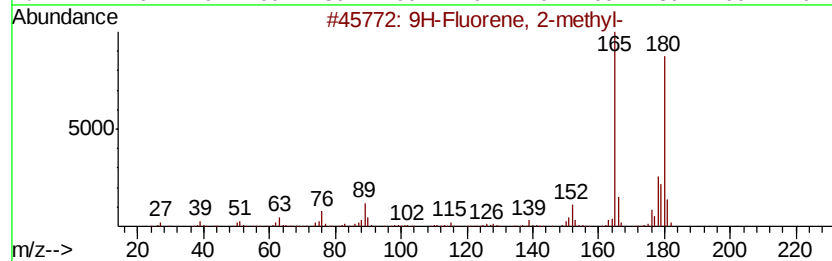
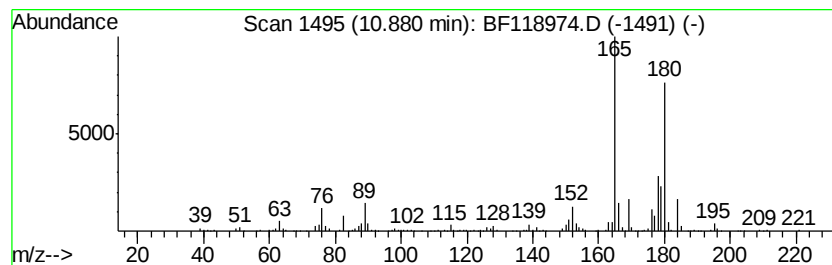
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 9H-Fluorene, 2-methyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.88	11.36 ng	709922	Phenanthrene-d10	11.27

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	9H-Fluorene, 2-methyl-	180	C14H12	001430-97-3	93
2		9H-Fluorene, 1-methyl-	180	C14H12	001730-37-6	93
3		4a,9a-Methano-9H-fluorene	180	C14H12	019540-84-2	86
4		9H-Fluorene, 3-methyl-	180	C14H12	002523-39-9	76
5		9H-Fluorene, 9-methyl-	180	C14H12	002523-37-7	74



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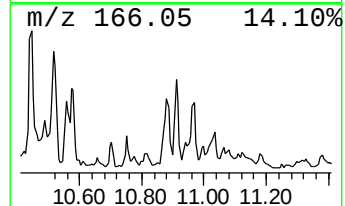
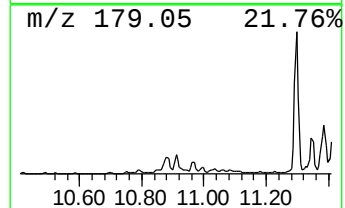
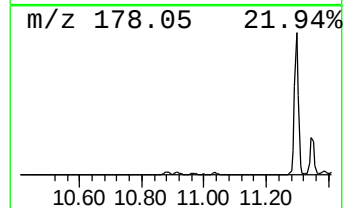
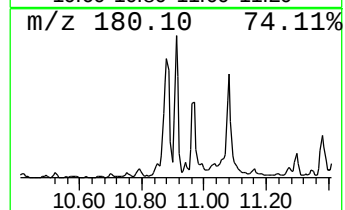
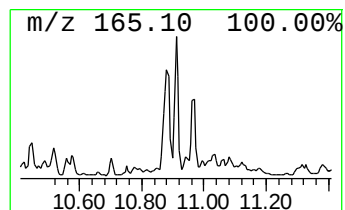
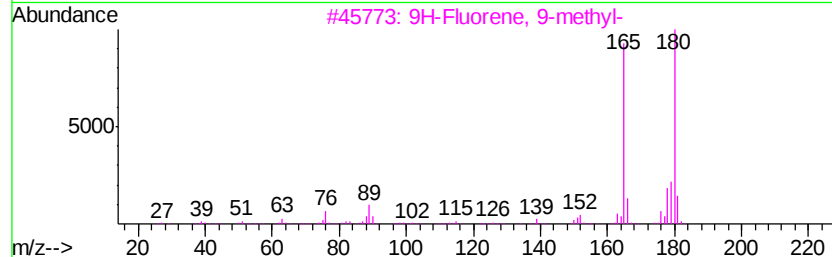
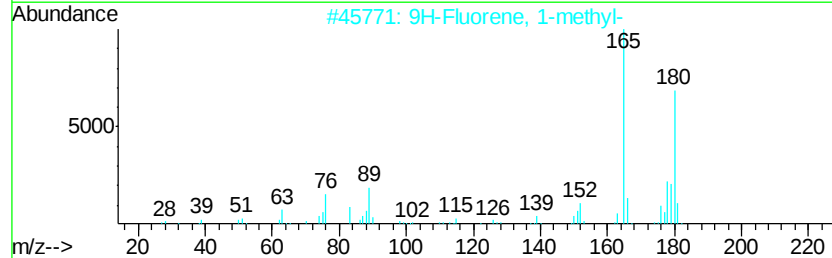
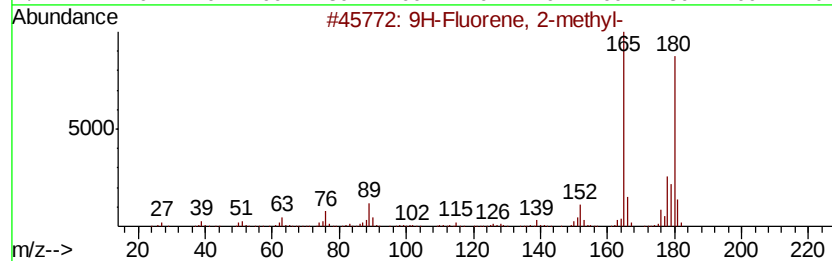
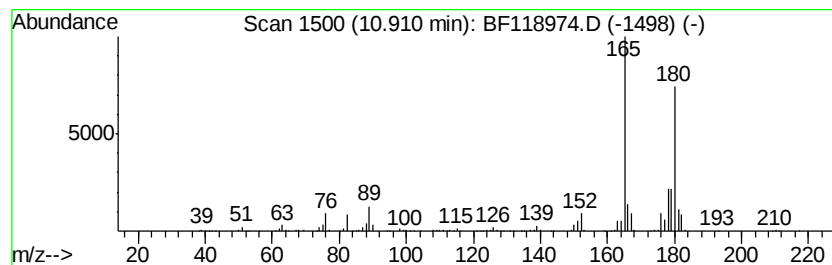
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 9H-Fluorene, 1-methyl- Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.91	7.64 ng	477519	Phenanthrene-d10	11.27

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		9H-Fluorene, 2-methyl-	180	C14H12	001430-97-3	97
2		9H-Fluorene, 1-methyl-	180	C14H12	001730-37-6	96
3		9H-Fluorene, 9-methyl-	180	C14H12	002523-37-7	94
4		9H-Fluorene, 3-methyl-	180	C14H12	002523-39-9	94
5		Ethylene, 1,1-diphenyl-	180	C14H12	000530-48-3	64



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF022020\
Data File : BF118974.D
Acq On : 20 Feb 2020 22:42
Operator : CG/JU
Sample : L1546-03
Misc :
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ClientSampleId :
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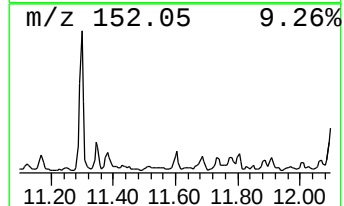
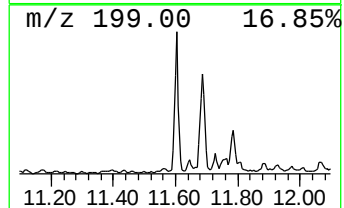
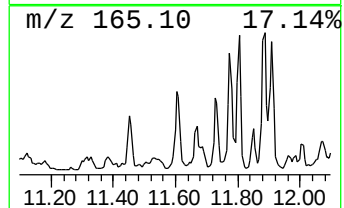
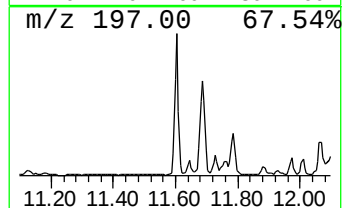
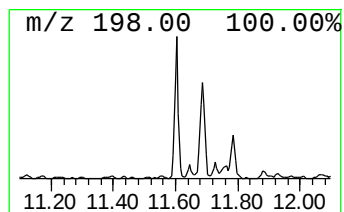
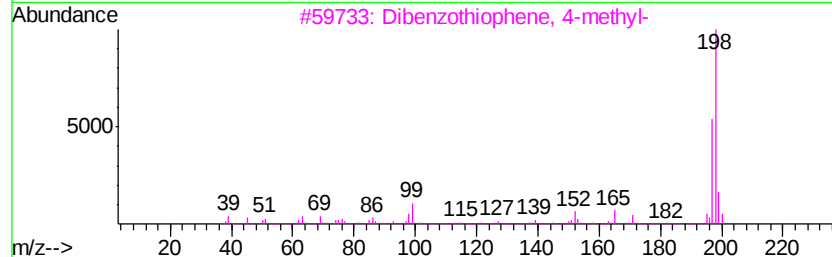
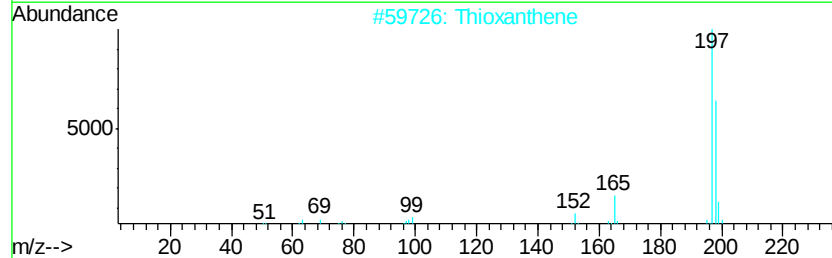
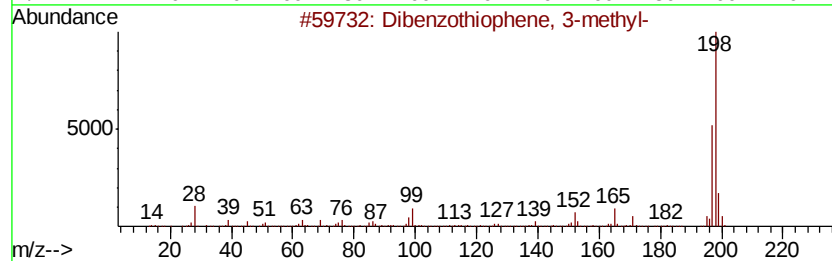
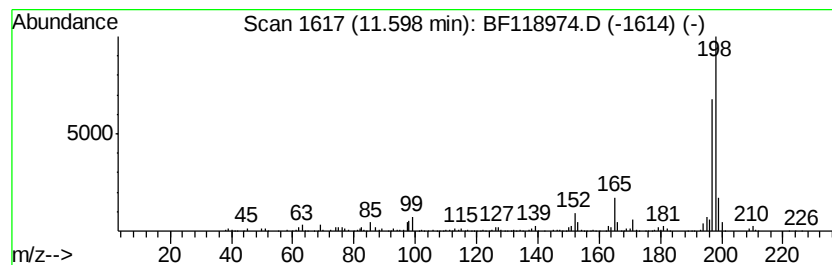
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Dibenzothiophene, 3-methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.60	11.74 ng	733812	Phenanthrene-d10	11.27

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Dibenzothiophene, 3-methyl-	198	C13H10S	016587-52-3	90
2		Thioxanthene	198	C13H10S	000261-31-4	86
3		Dibenzothiophene, 4-methyl-	198	C13H10S	007372-88-5	81
4		Tacrine	198	C13H14N2	000321-64-2	64
5		Benzenamine, 4,4'-methylenebis-	198	C13H14N2	000101-77-9	59



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF022020\
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Acq On : 20 Feb 2020 22:42
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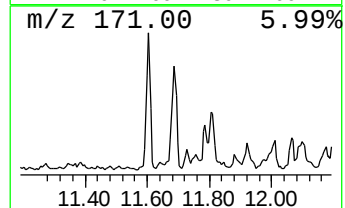
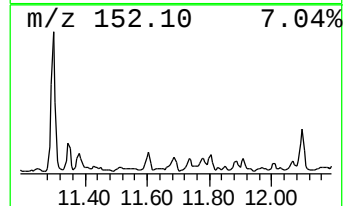
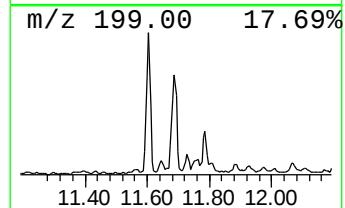
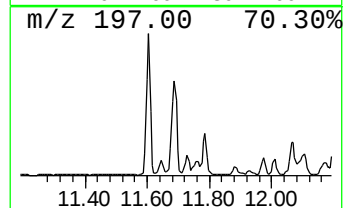
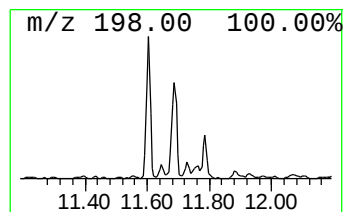
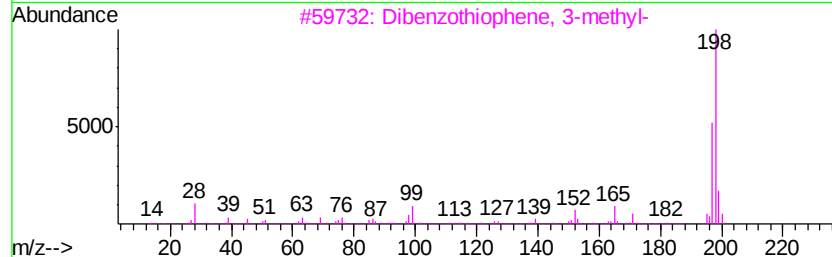
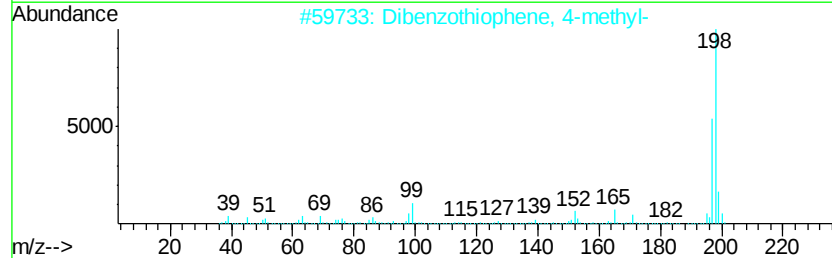
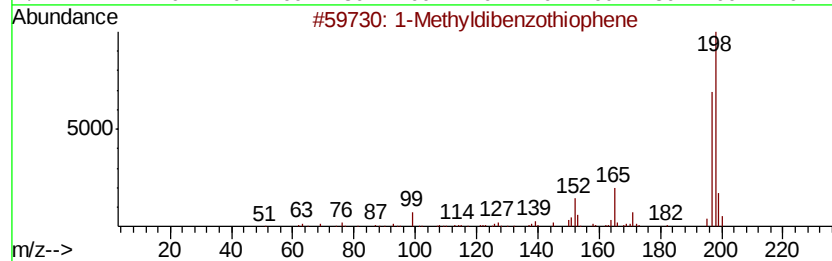
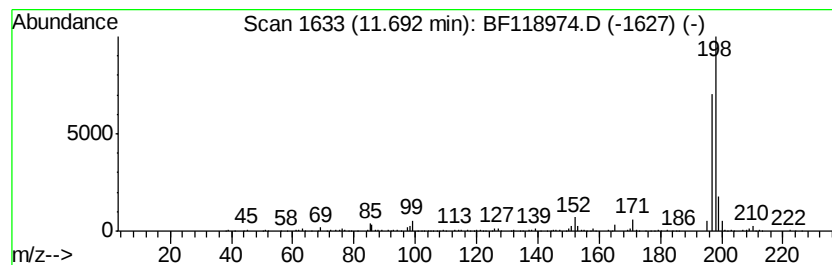
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 1-Methyldibenzothiophene Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.69	9.35 ng	584546	Phenanthrene-d10	11.27

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Methyldibenzothiophene	198	C13H10S	031317-07-4	91
2		Dibenzothiophene, 4-methyl-	198	C13H10S	007372-88-5	90
3		Dibenzothiophene, 3-methyl-	198	C13H10S	016587-52-3	90
4		4-Methylnaphtho[1,2-b]thiophene	198	C13H10S	067388-11-8	90
5		Tacrine	198	C13H14N2	000321-64-2	80



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF022020\
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Acq On : 20 Feb 2020 22:42
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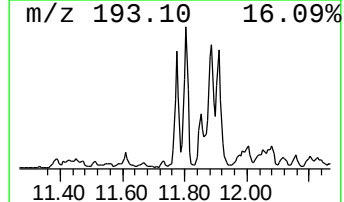
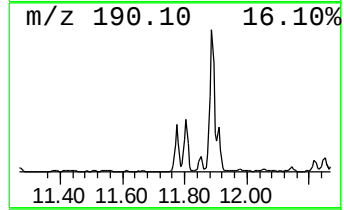
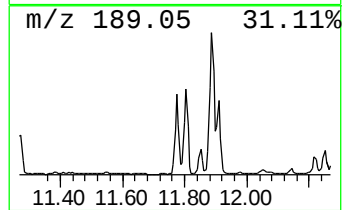
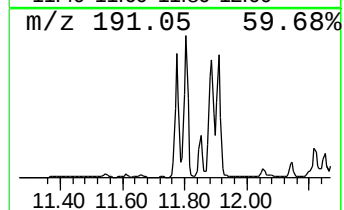
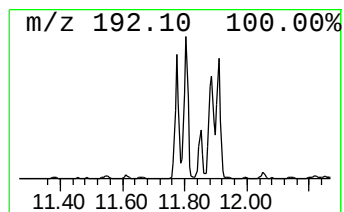
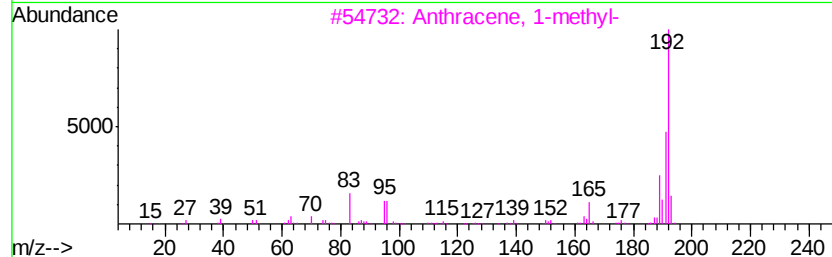
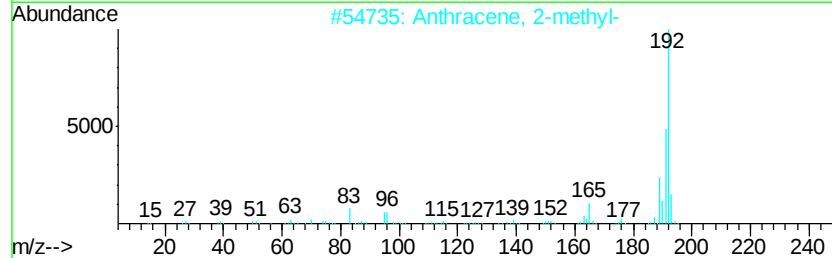
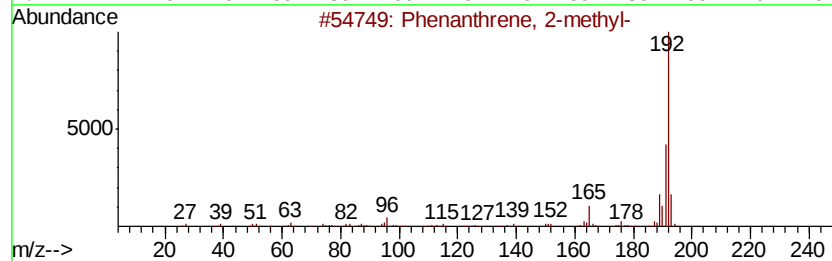
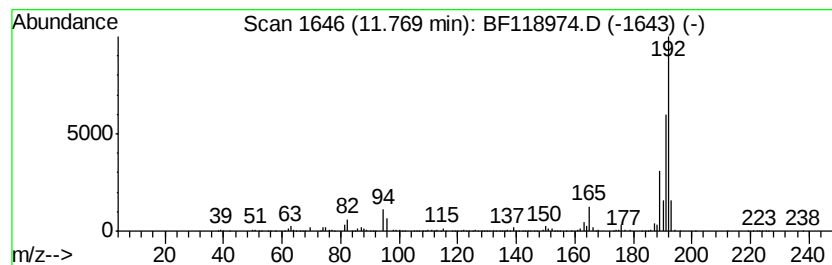
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Phenanthrene, 2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.77	25.47 ng	1591590	Phenanthrene-d10	11.27

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	95
2		Anthracene, 2-methyl-	192	C15H12	000613-12-7	95
3		Anthracene, 1-methyl-	192	C15H12	000610-48-0	94
4		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	91
5		Anthracene, 9-methyl-	192	C15H12	000779-02-2	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF022020\
Data File : BF118974.D
Acq On : 20 Feb 2020 22:42
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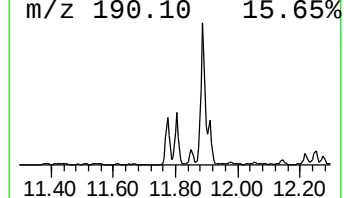
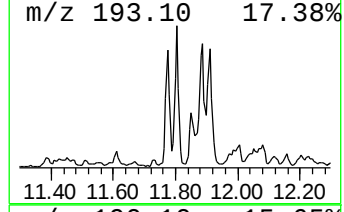
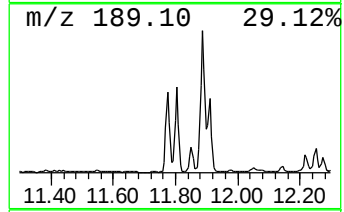
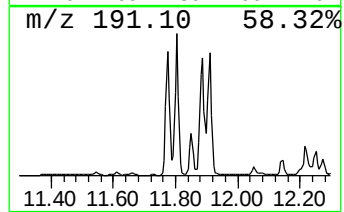
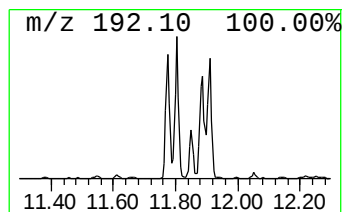
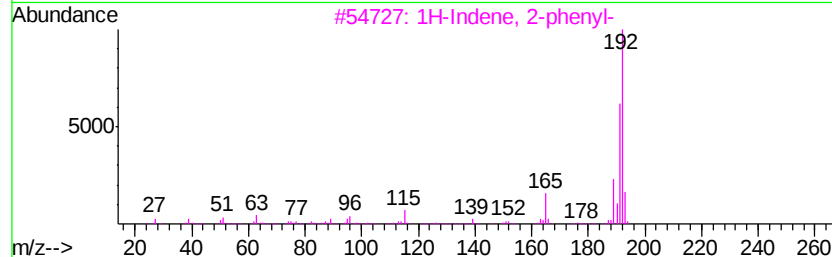
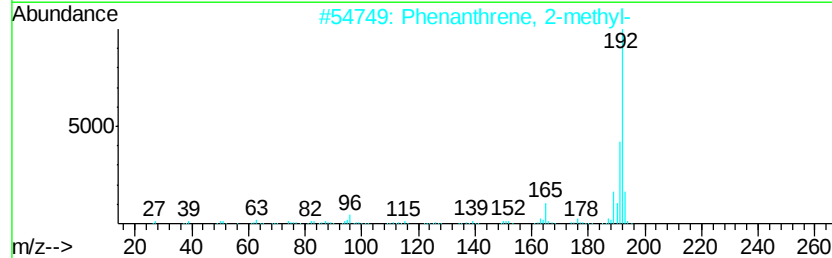
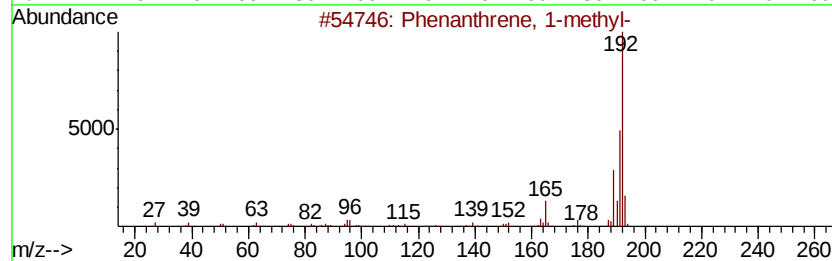
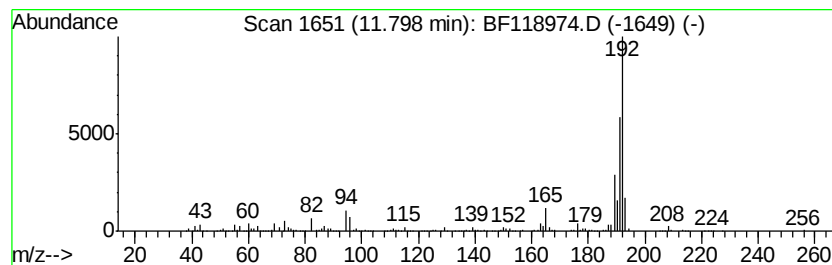
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Phenanthrene, 1-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.80	33.55 ng	2096750	Phenanthrene-d10	11.27

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	96
2		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	95
3		1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	95
4		Anthracene, 2-methyl-	192	C15H12	000613-12-7	95
5		1H-Cyclopropa[1]phenanthrene, 1a,...	192	C15H12	000949-41-7	93



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF022020\
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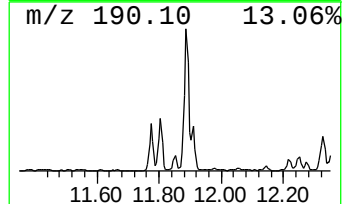
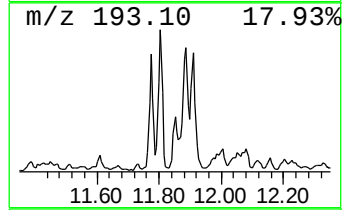
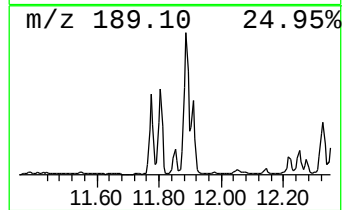
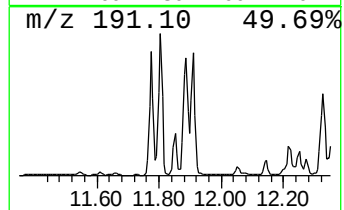
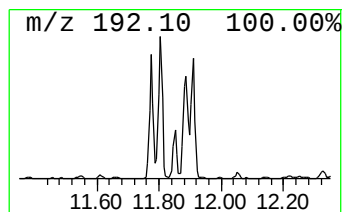
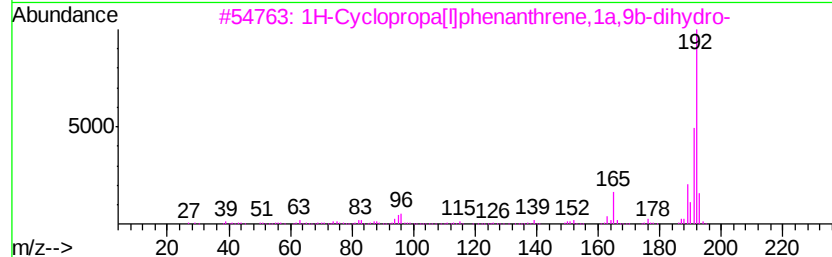
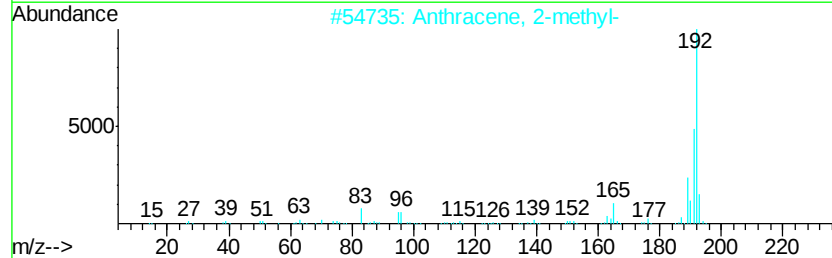
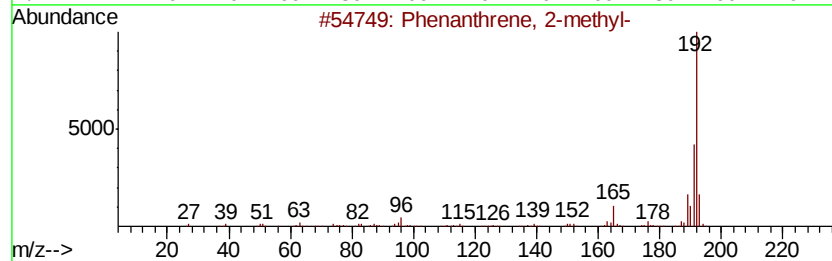
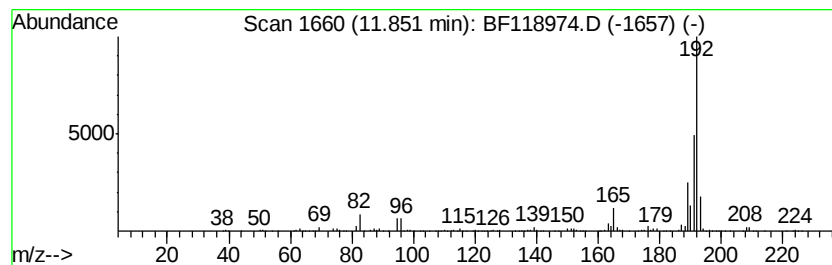
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Anthracene, 2-methyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.85	8.95 ng	559510	Phenanthrene-d10	11.27	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	97
2	Anthracene, 2-methyl-	192	C15H12	000613-12-7	96
3	1H-Cyclopropa[1]phenanthrene,1a,...	192	C15H12	000949-41-7	96
4	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	96
5	Phenanthrene, 3-methyl-	192	C15H12	000832-71-3	95



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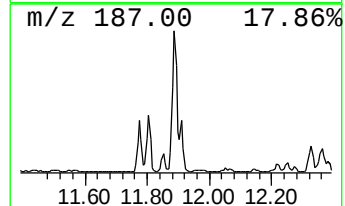
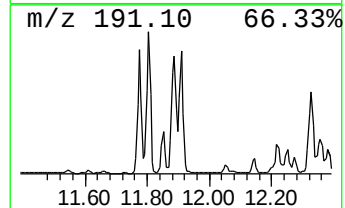
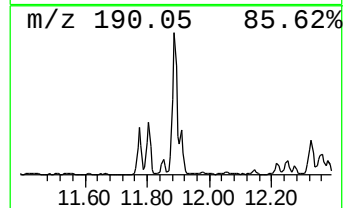
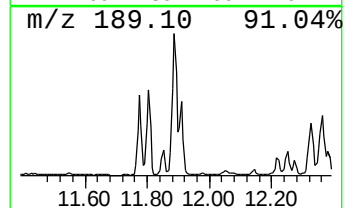
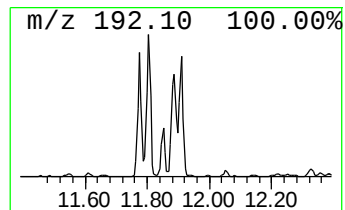
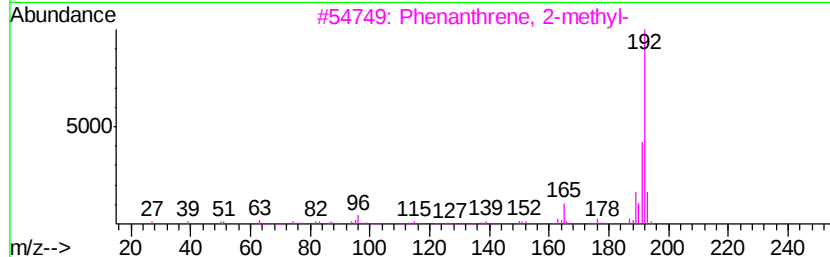
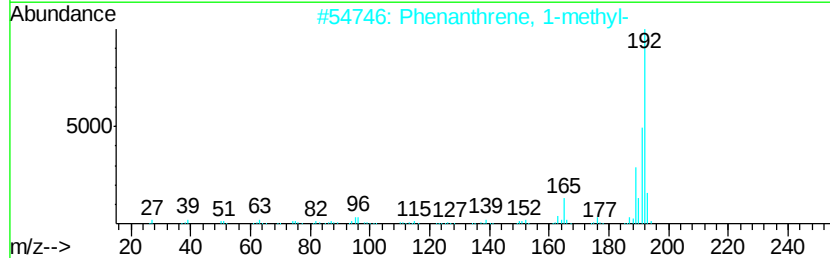
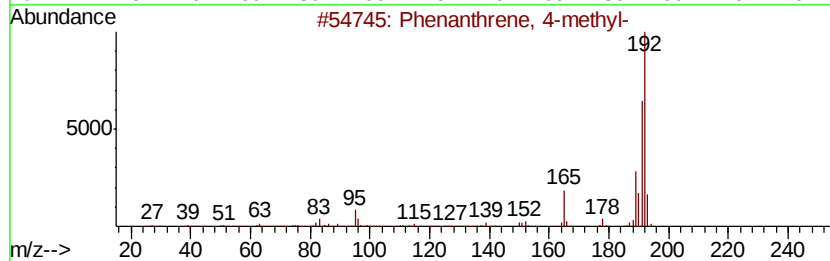
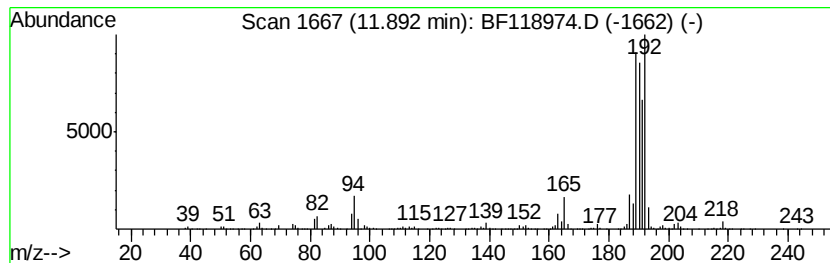
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Phenanthrene, 4-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.89	35.57 ng	2223030	Phenanthrene-d10	11.27	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	90
2	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	60
3	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	60
4	Anthracene, 1-methyl-	192	C15H12	000610-48-0	60
5	Anthracene, 2-methyl-	192	C15H12	000613-12-7	60



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Data File : BF118974.D
Acq On : 20 Feb 2020 22:42
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ALS Vial : 10 Sample Multiplier: 1

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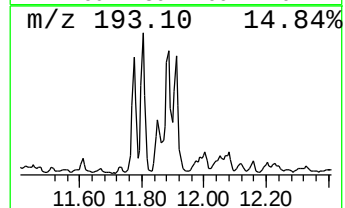
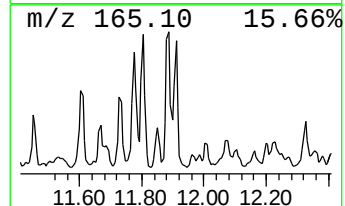
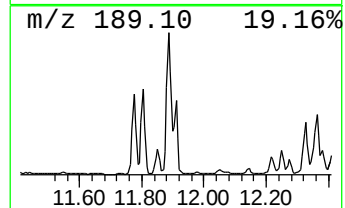
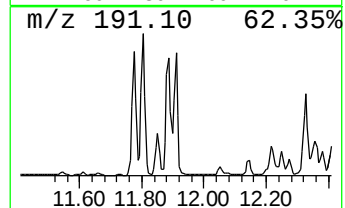
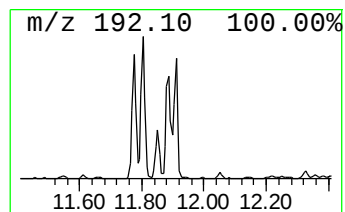
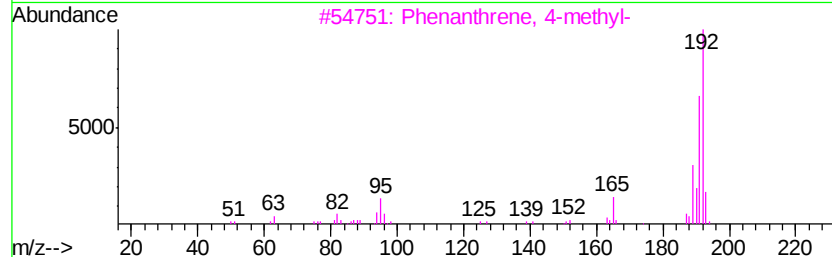
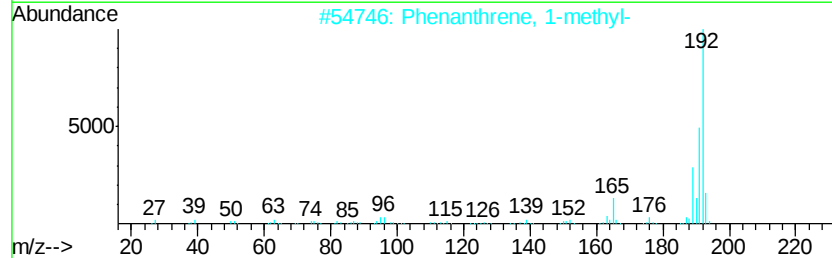
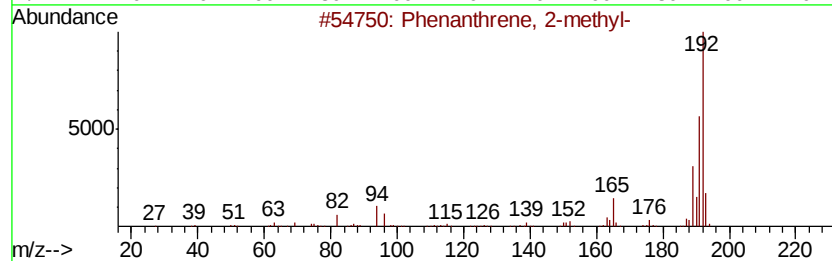
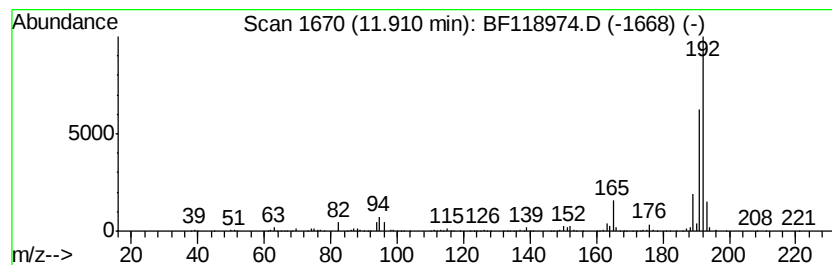
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 1H-Cyclopropa[1]phenanthren... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.91	22.34 ng	1396220	Phenanthrene-d10	11.27

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	91
2			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	90
3			Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	90
4			1H-Cyclopropa[1]phenanthrene, 1a, ...	192	C15H12	000949-41-7	90
5			Anthracene, 2-methyl-	192	C15H12	000613-12-7	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF022020\
Data File : BF118974.D
Acq On : 20 Feb 2020 22:42
Operator : CG/JU
Sample : L1546-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

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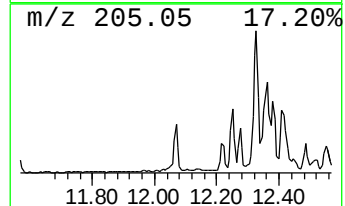
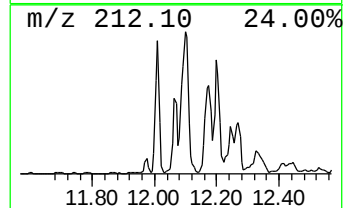
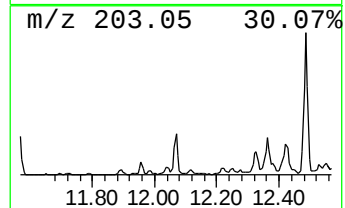
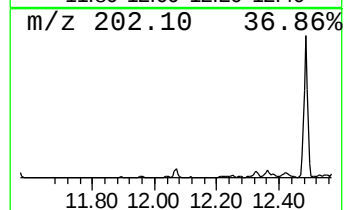
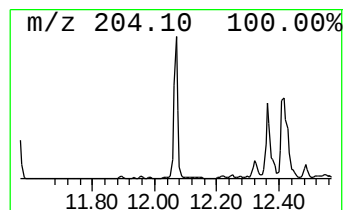
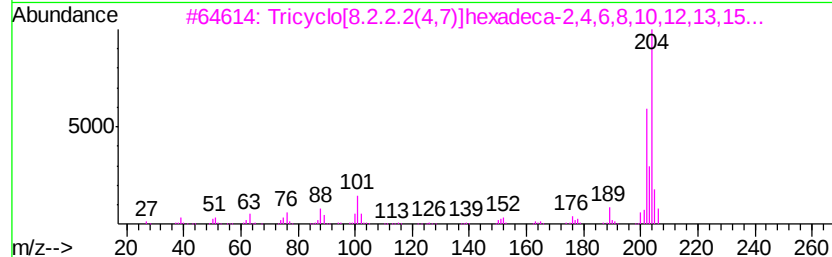
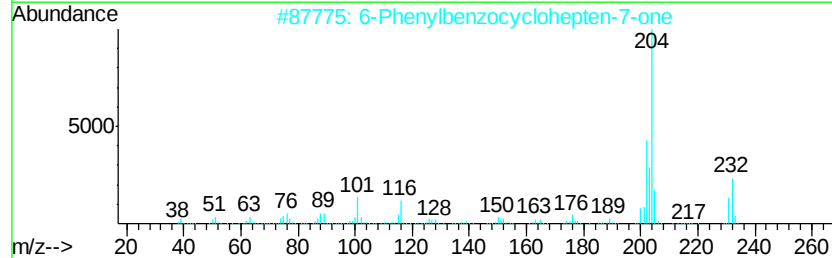
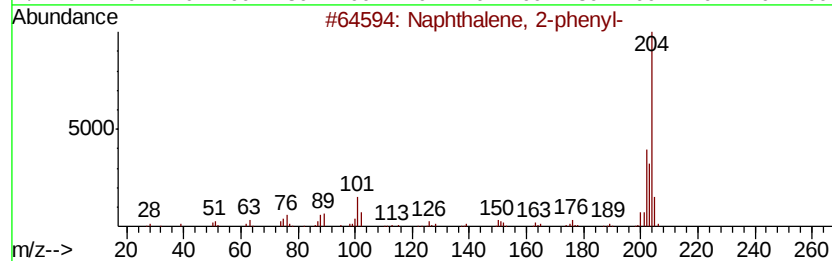
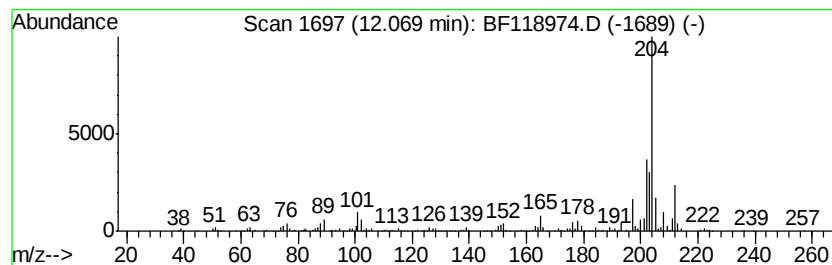
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Naphthalene, 2-phenyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.07	15.65 ng	977942	Phenanthrene-d10	11.27

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	93
2		6-Phenylbenzocyclohepten-7-one	232	C17H12O	093327-56-1	64
3		Tricyclo[8.2.2.2(4,7)]hexadeca-2...	204	C16H12	006572-60-7	64
4		6-Cyanophenanthridine	204	C14H8N2	002176-28-5	50
5		Benzene, 1,1'-(1-buten-3-yne-1,4...	204	C16H12	013141-45-2	49



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF022020\
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Acq On : 20 Feb 2020 22:42
Operator : CG/JU
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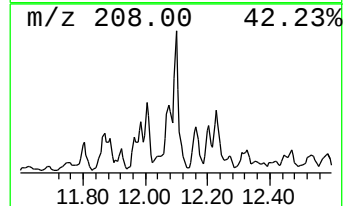
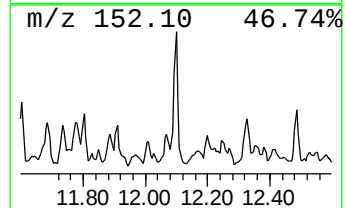
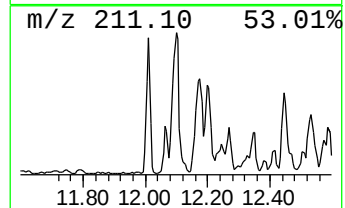
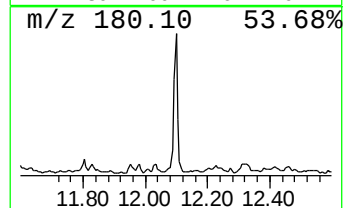
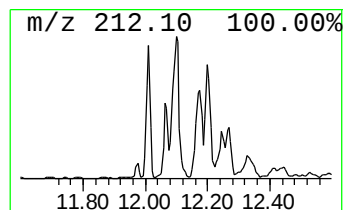
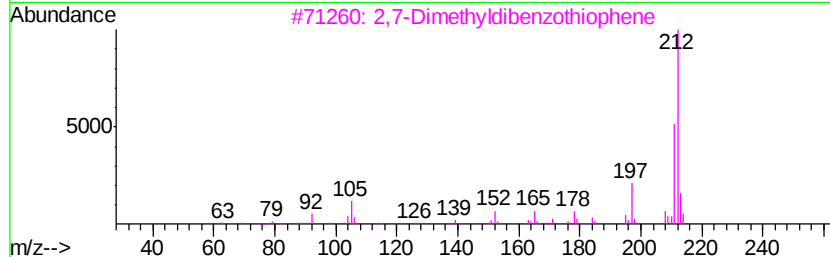
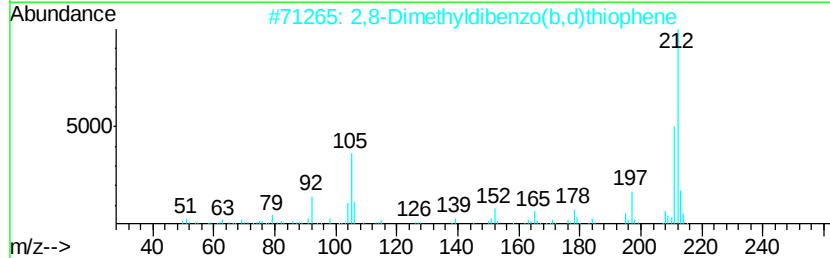
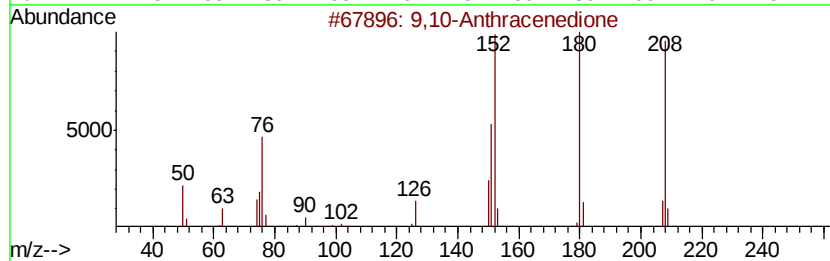
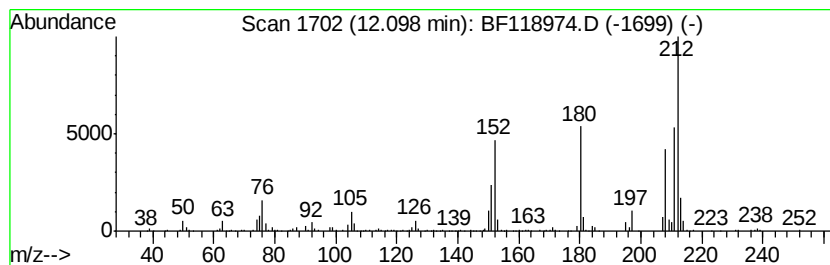
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 9,10-Anthracenedione Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.10	11.69 ng	730497	Phenanthrene-d10	11.27

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			9,10-Anthracenedione	208	C14H8O2	000084-65-1	89
2			2,8-Dimethyldibenzo(b,d)thiophene	212	C14H12S	001207-15-4	86
3			2,7-Dimethyldibenzothiophene	212	C14H12S	031317-19-8	43
4			3,7-Dimethyldibenzothiophene	212	C14H12S	001136-85-2	43
5			1,7-Dimethyldibenzothiophene	212	C14H12S	089816-53-5	43



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF022020\
Data File : BF118974.D
Acq On : 20 Feb 2020 22:42
Operator : CG/JU
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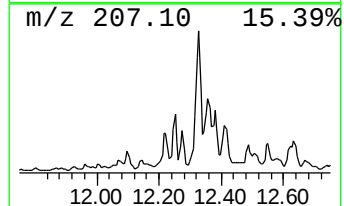
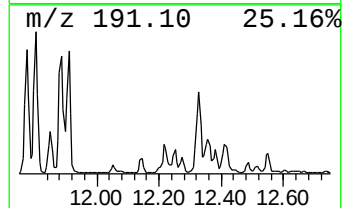
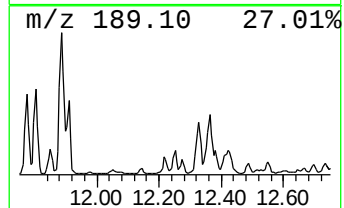
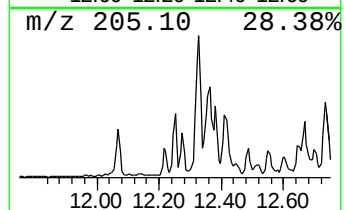
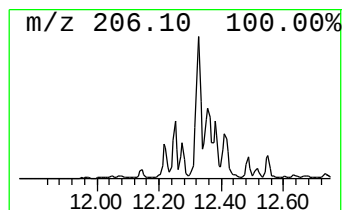
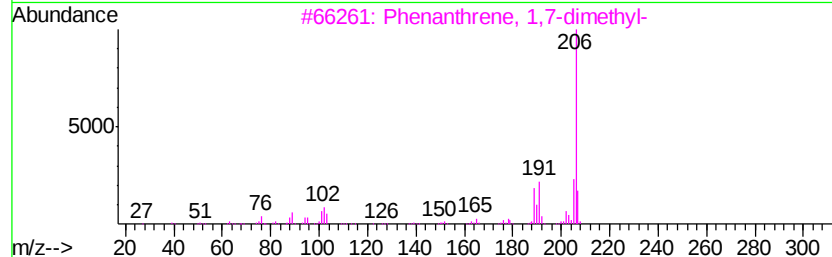
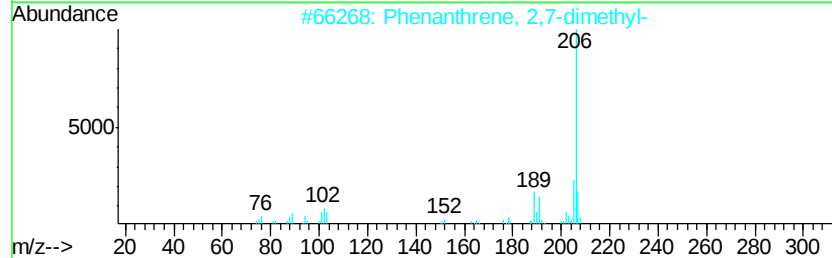
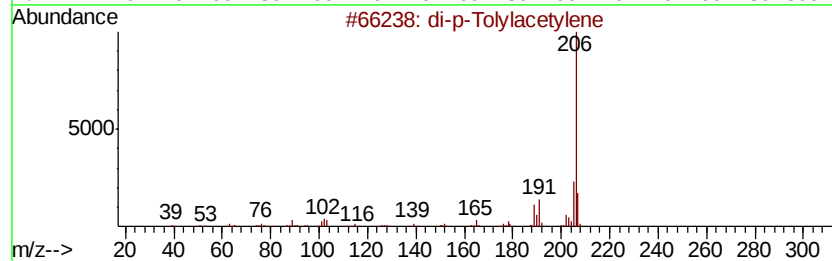
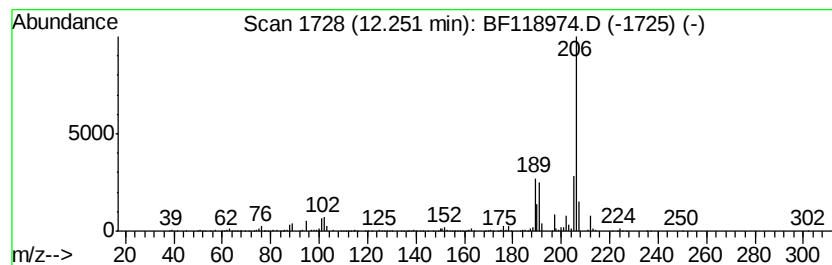
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 14 di-p-Tolylacetylene Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.25	7.82 ng	488569	Phenanthrene-d10	11.27

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			di-p-Tolylacetylene	206	C16H14	002789-88-0	90
2			Phenanthrene, 2,7-dimethyl-	206	C16H14	001576-69-8	87
3			Phenanthrene, 1,7-dimethyl-	206	C16H14	000483-87-4	86
4			1,4-Diphenyl-1,3-butadiene	206	C16H14	000886-65-7	80
5			Phenanthrene, 3,6-dimethyl-	206	C16H14	001576-67-6	80



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF022020\
Data File : BF118974.D
Acq On : 20 Feb 2020 22:42
Operator : CG/JU
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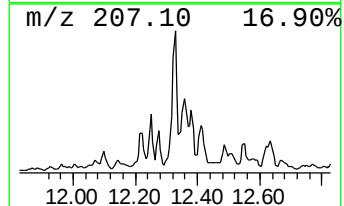
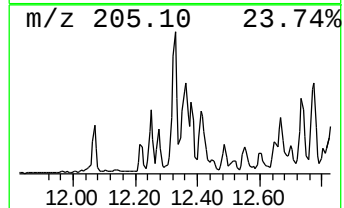
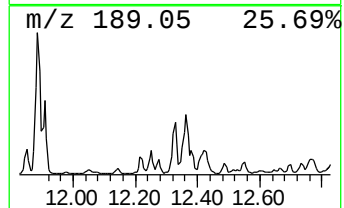
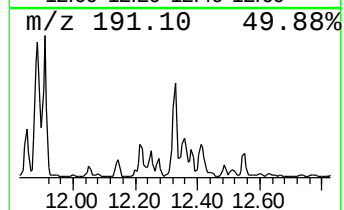
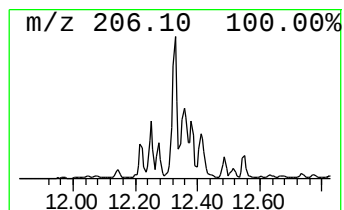
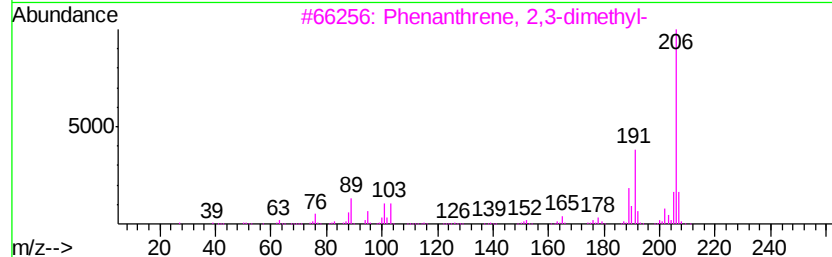
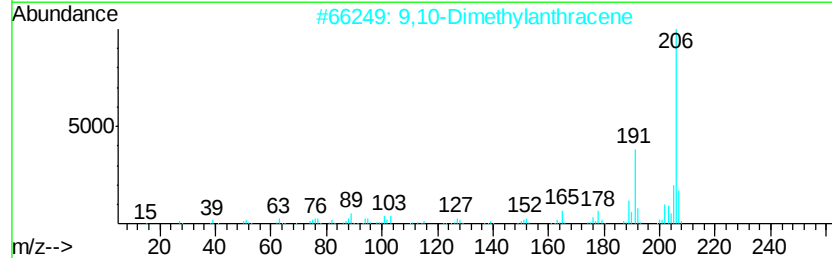
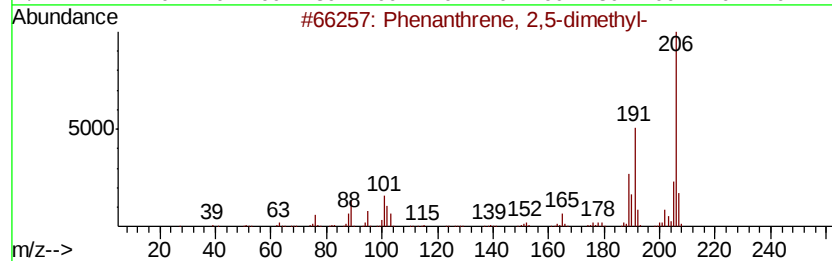
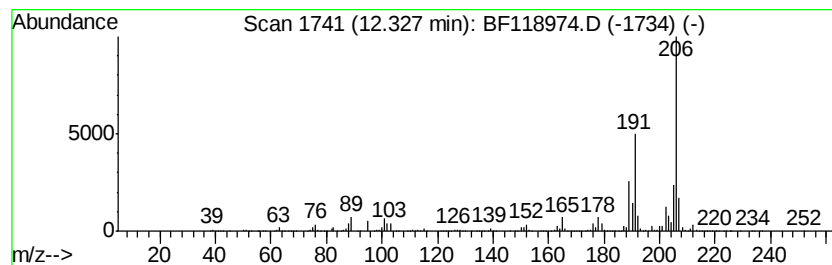
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 15 Phenanthrene, 2,5-dimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.33	29.03 ng	1813850	Phenanthrene-d10	11.27

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	96
2		9,10-Dimethylantracene	206	C16H14	000781-43-1	95
3		Phenanthrene, 2,3-dimethyl-	206	C16H14	003674-65-5	93
4		Phenanthrene, 1,7-dimethyl-	206	C16H14	000483-87-4	90
5		Anthracene, 1,4-dimethyl-	206	C16H14	000781-92-0	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF022020\
Data File : BF118974.D
Acq On : 20 Feb 2020 22:42
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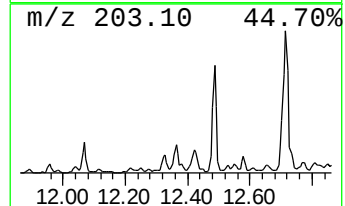
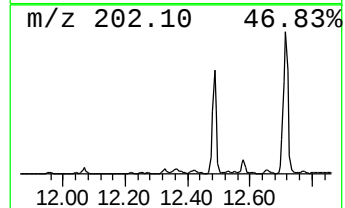
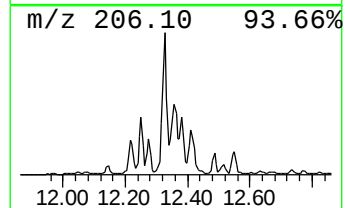
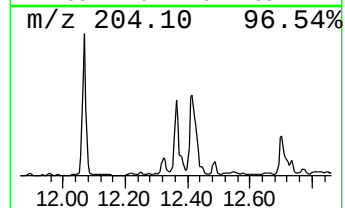
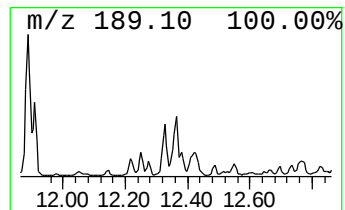
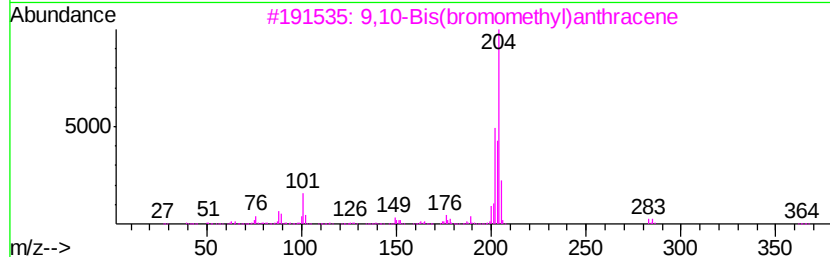
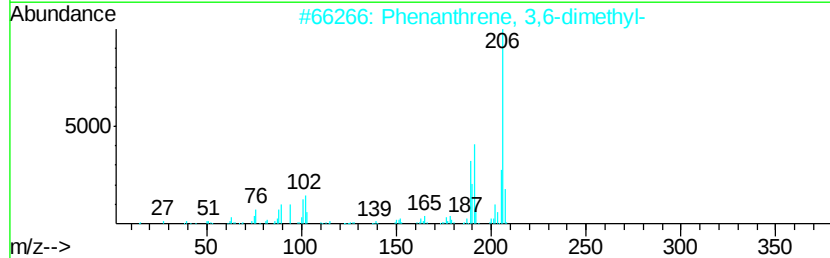
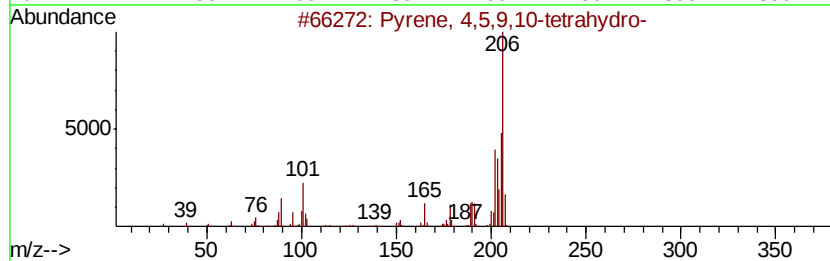
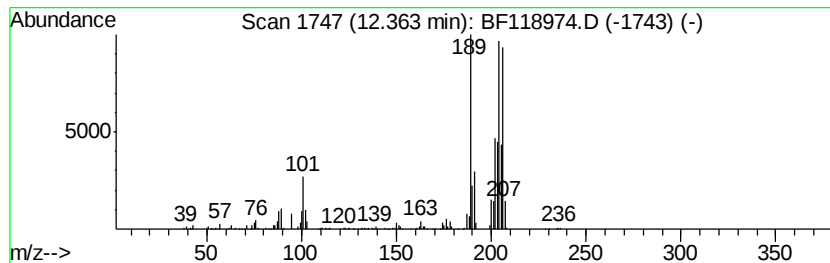
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 unknown12.36 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.36	19.00 ng	1187040	Phenanthrene-d10	11.27

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pyrene, 4,5,9,10-tetrahydro-	206	C16H14	000781-17-9	46
2		Phenanthrene, 3,6-dimethyl-	206	C16H14	001576-67-6	38
3		9,10-Bis(bromomethyl)anthracene	362	C16H12Br2	034373-96-1	30
4		Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	25
5		Tetramisole	204	C11H12N2S	005036-02-2	25



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF022020\
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Acq On : 20 Feb 2020 22:42
Operator : CG/JU
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ClientSampleId :
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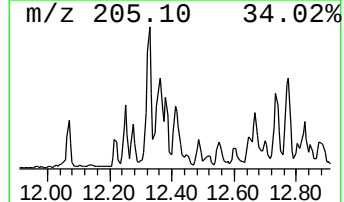
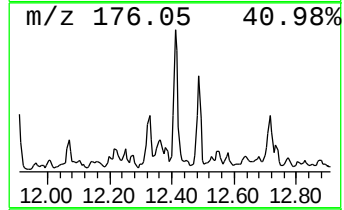
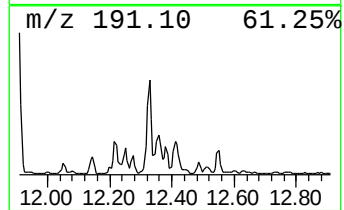
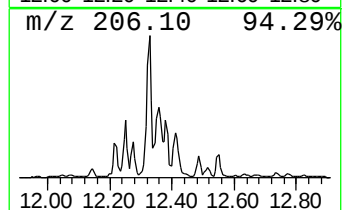
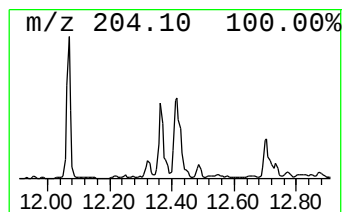
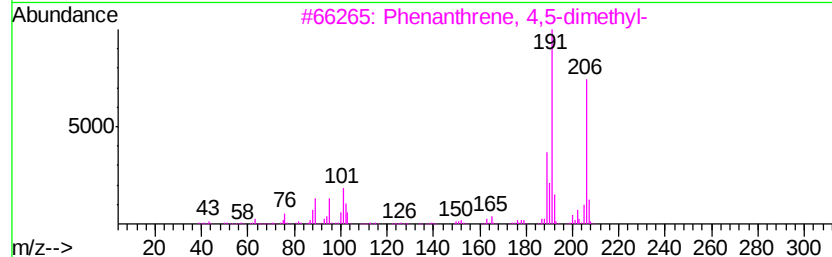
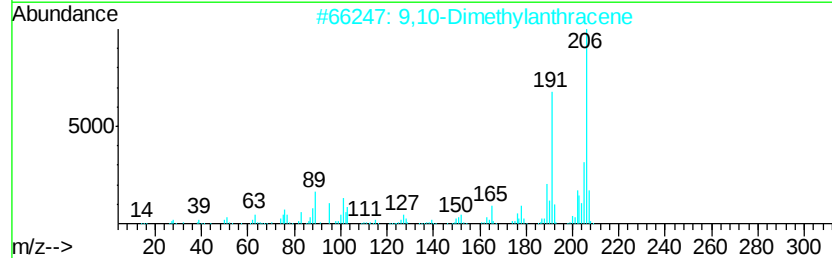
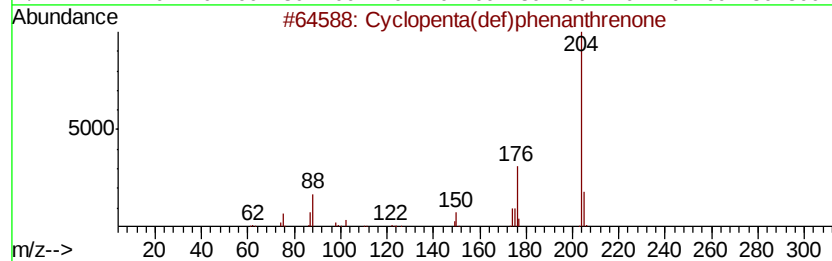
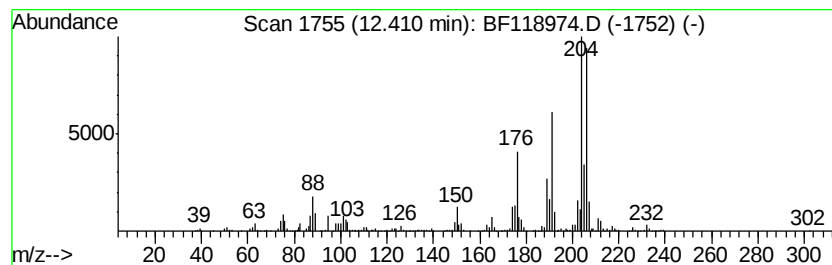
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 17 Cyclopenta(def)phenanthrenone Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.41	15.20 ng	949878	Phenanthrene-d10	11.27

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopenta(def)phenanthrenone	204	C15H8O	005737-13-3	93
2		9,10-Dimethylantracene	206	C16H14	000781-43-1	70
3		Phenanthrene, 4,5-dimethyl-	206	C16H14	003674-69-9	64
4		1H-Indene, 2-methyl-3-phenyl-	206	C16H14	035099-60-6	55
5		1,3-Diphenyl-3-methylcyclopropene	206	C16H14	086544-79-8	55



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF022020\
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Misc :
ALS Vial : 10 Sample Multiplier: 1

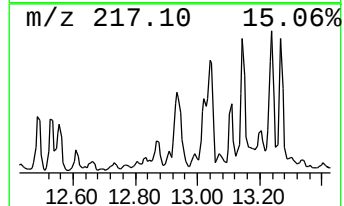
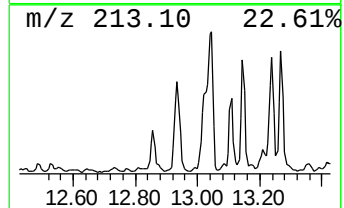
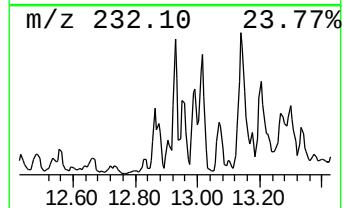
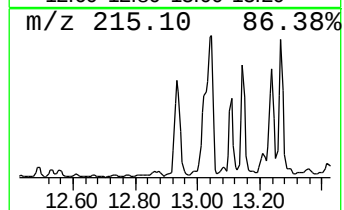
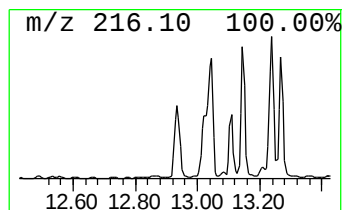
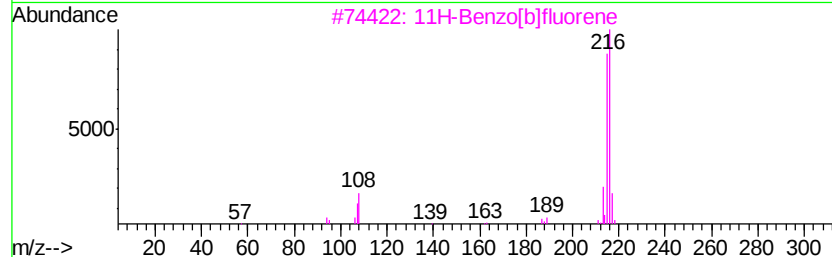
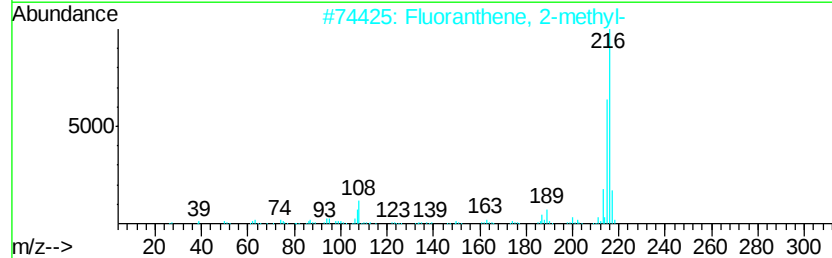
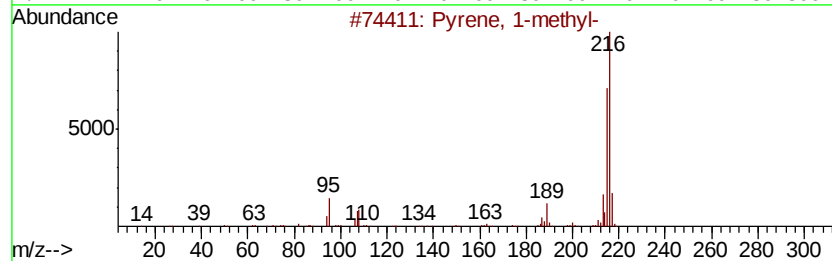
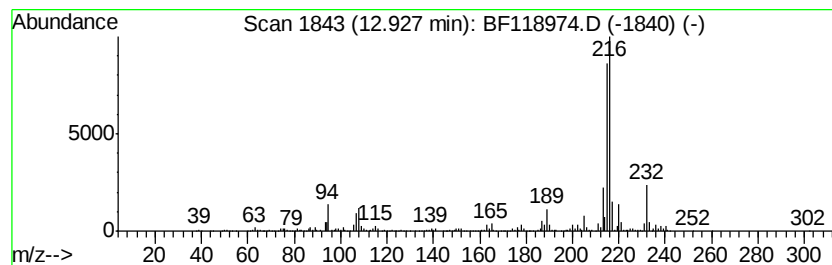
Instrument :
BNA_F
ClientSampleId :
INDIAN-ISLAND-(2)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 11H-Benzo[b]fluorene Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.93	7.94 ng	1271440	Chrysene-d12	13.91	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pyrene, 1-methyl-	216	C17H12	002381-21-7	96
2	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	90
3	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	89
4	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	80
5	7H-Benzo[c]fluorene	216	C17H12	000205-12-9	68



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF022020\
Data File : BF118974.D
Acq On : 20 Feb 2020 22:42
Operator : CG/JU
Sample : L1546-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
INDIAN-ISLAND-(2)

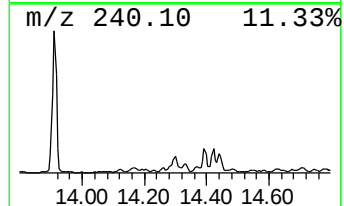
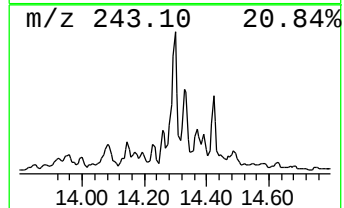
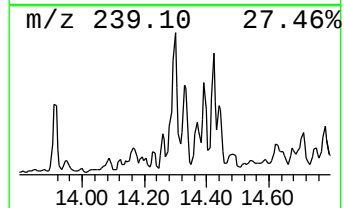
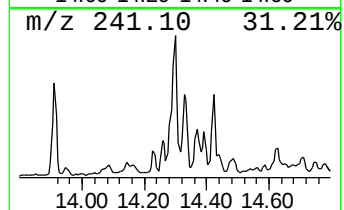
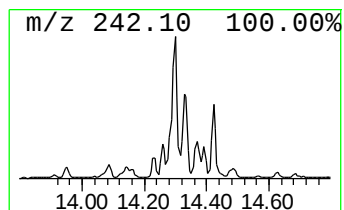
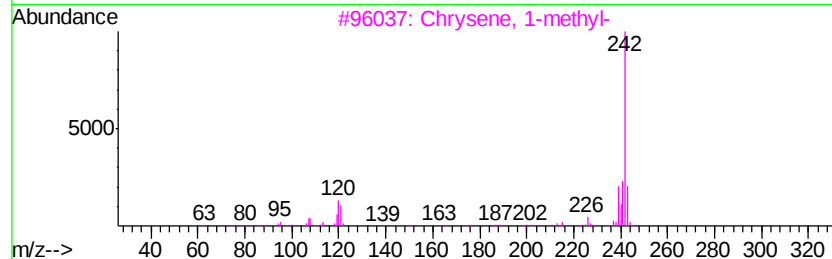
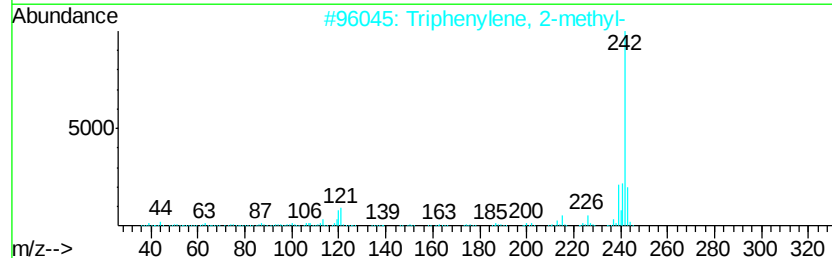
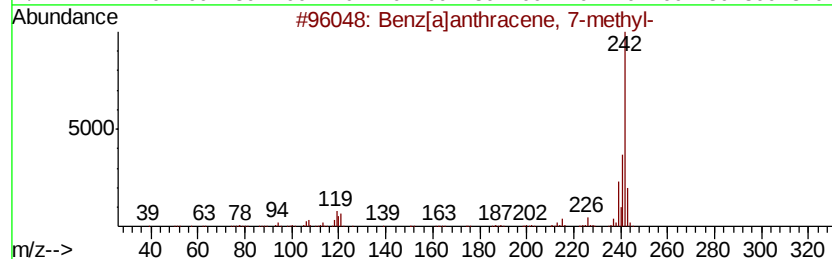
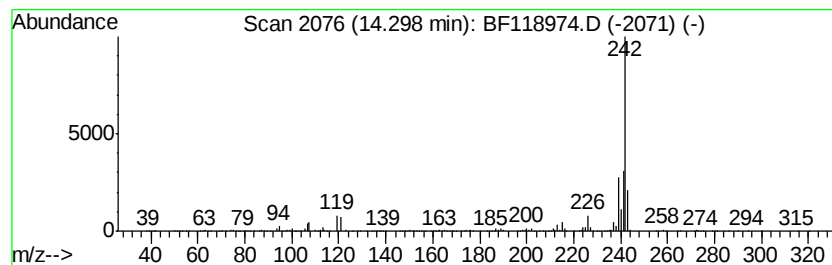
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 19 Benz[a]anthracene, 7-methyl- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.30	8.86 ng	1418030	Chrysene-d12	13.91

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benz[a]anthracene, 7-methyl-	242	C19H14	002541-69-7	93
2			Triphenylene, 2-methyl-	242	C19H14	001705-84-6	93
3			Chrysene, 1-methyl-	242	C19H14	003351-28-8	91
4			Chrysene, 5-methyl-	242	C19H14	003697-24-3	90
5			Chrysene, 3-methyl-	242	C19H14	003351-31-3	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF022020\
Data File : BF118974.D
Acq On : 20 Feb 2020 22:42
Operator : CG/JU
Sample : L1546-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
INDIAN-ISLAND-(2)

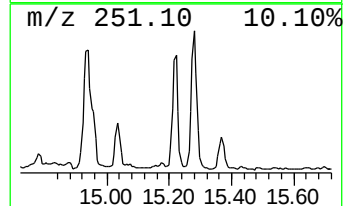
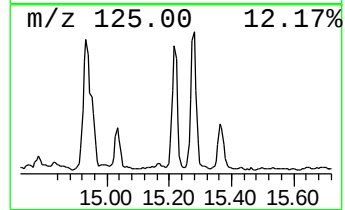
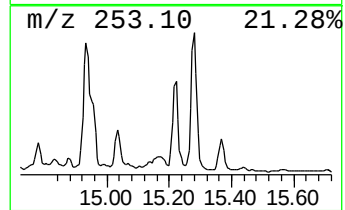
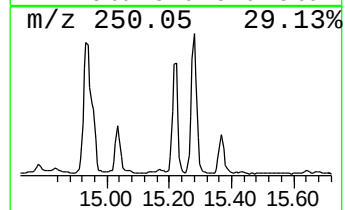
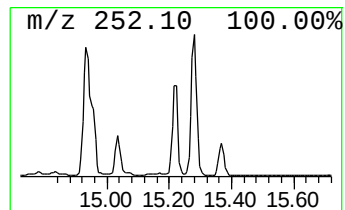
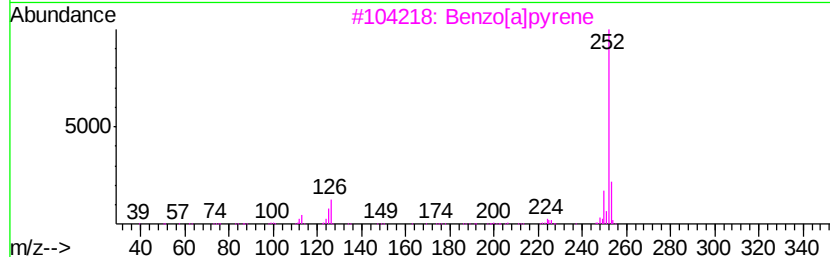
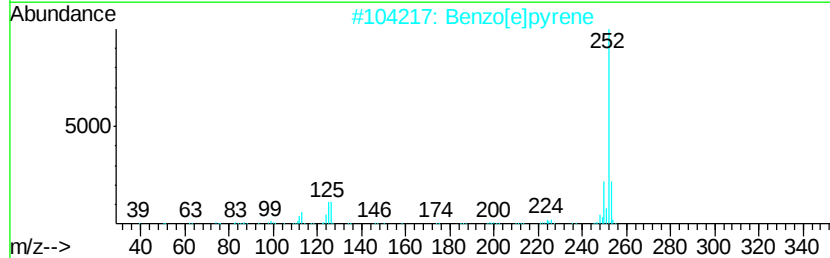
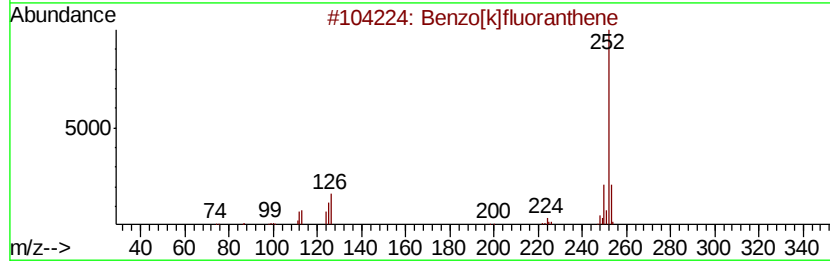
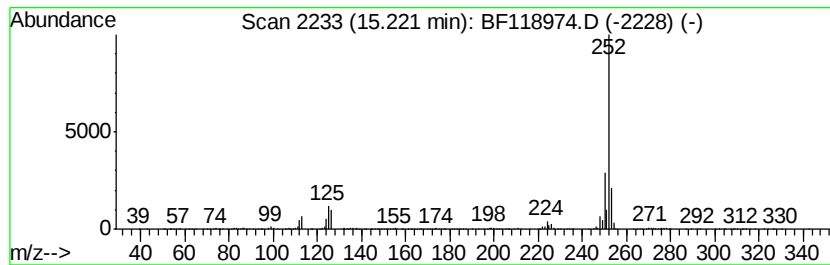
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 20 Benzo[e]pyrene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.22	17.81 ng	976995	Perylene-d12	15.34

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzo[k]fluoranthene	252	C20H12	000207-08-9	98
2		Benzo[e]pyrene	252	C20H12	000192-97-2	98
3		Benzo[a]pyrene	252	C20H12	000050-32-8	97
4		Benz[el]acephenanthrylene	252	C20H12	000205-99-2	95
5		Perylene	252	C20H12	000198-55-0	94



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF022020\
 Data File : BF118974.D
 Acq On : 20 Feb 2020 22:42
 Operator : CG/JU
 Sample : L1546-03
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 INDIAN-ISLAND-(2)

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF022020.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
Naphthalene, 1,5-...	9.45	9.0 ng		747692	3	9.79	1668230	20.0
9H-Fluorene, 2-me...	10.88	11.4 ng		709922	4	11.27	1249800	20.0
9H-Fluorene, 1-me...	10.91	7.6 ng		477519	4	11.27	1249800	20.0
Dibenzothiophene,...	11.60	11.7 ng		733812	4	11.27	1249800	20.0
1-Methyldibenzoth...	11.69	9.3 ng		584546	4	11.27	1249800	20.0
Phenanthrene, 2-m...	11.77	25.5 ng		1591590	4	11.27	1249800	20.0
Phenanthrene, 1-m...	11.80	33.5 ng		2096750	4	11.27	1249800	20.0
Anthracene, 2-met...	11.85	8.9 ng		559510	4	11.27	1249800	20.0
Phenanthrene, 4-m...	11.89	35.6 ng		2223030	4	11.27	1249800	20.0
1H-Cyclopropa[1]p...	11.91	22.3 ng		1396220	4	11.27	1249800	20.0
Naphthalene, 2-ph...	12.07	15.7 ng		977942	4	11.27	1249800	20.0
9,10-Anthracenedione	12.10	11.7 ng		730497	4	11.27	1249800	20.0
di-p-Tolylacetylene	12.25	7.8 ng		488569	4	11.27	1249800	20.0
Phenanthrene, 2,5...	12.33	29.0 ng		1813850	4	11.27	1249800	20.0
unknown12.36	12.36	19.0 ng		1187040	4	11.27	1249800	20.0
Cyclopenta(def)ph...	12.41	15.2 ng		949878	4	11.27	1249800	20.0
11H-Benzo[b]fluorene	12.93	7.9 ng		1271440	5	13.91	3202470	20.0
Benz[a]anthracene...	14.30	8.9 ng		1418030	5	13.91	3202470	20.0
Benzo[e]pyrene	15.22	17.8 ng		976995	6	15.34	1097360	20.0