

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF022520\
 Data File : BF119085.D
 Acq On : 26 Feb 2020 1:10
 Operator : CG/JU
 Sample : L1367-05 5X
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :

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	5.375	556	559	579	rBV	604618	651198	16.96%	1.253%
2	6.392	729	732	748	rBV	653200	664260	17.30%	1.278%
3	6.751	789	793	802	rBV	1132430	999480	26.03%	1.923%
4	6.910	816	820	828	rBV	689972	635846	16.56%	1.223%
5	7.310	885	888	895	rBV	439483	407178	10.60%	0.783%
6	8.033	1006	1011	1014	rBV	1492970	1272733	33.15%	2.448%
7	8.745	1127	1132	1136	rBV	145521	123454	3.22%	0.237%
8	8.845	1143	1149	1158	rVB	184060	172586	4.49%	0.332%
9	9.110	1190	1194	1205	rVB	1003891	958375	24.96%	1.844%
10	9.374	1235	1239	1243	rBV	122043	165237	4.30%	0.318%
11	9.445	1247	1251	1254	rVV	232164	236697	6.16%	0.455%
12	9.474	1254	1256	1258	rVV	135961	114911	2.99%	0.221%
13	9.510	1258	1262	1266	rVV2	93073	120910	3.15%	0.233%
14	9.563	1268	1271	1276	rVV	116506	128809	3.35%	0.248%
15	9.645	1280	1285	1290	rBV	402705	360124	9.38%	0.693%
16	9.786	1302	1309	1312	rBV	1598945	1425553	37.13%	2.742%
17	9.816	1312	1314	1318	rVB	322233	273655	7.13%	0.526%
18	9.986	1340	1343	1347	rVV	183275	192241	5.01%	0.370%
19	10.027	1347	1350	1353	rVB	104241	85657	2.23%	0.165%
20	10.104	1358	1363	1365	rBV2	100020	123315	3.21%	0.237%
21	10.133	1365	1368	1374	rVB	73333	85283	2.22%	0.164%
22	10.192	1374	1378	1380	rBV2	65127	89564	2.33%	0.172%
23	10.333	1399	1402	1408	rVB	348875	363355	9.46%	0.699%
24	10.398	1408	1413	1415	rBV2	80527	103017	2.68%	0.198%
25	10.445	1418	1421	1425	rVV4	53587	67028	1.75%	0.129%
26	10.486	1425	1428	1432	rVB	77945	80182	2.09%	0.154%
27	10.574	1437	1443	1447	rBV	528899	567268	14.77%	1.091%
28	10.698	1460	1464	1470	rVB3	119048	143043	3.73%	0.275%
29	10.880	1490	1495	1497	rBV3	114085	134604	3.51%	0.259%
30	10.910	1497	1500	1503	rVV	87148	84545	2.20%	0.163%
31	11.010	1512	1517	1519	rBV4	58792	82721	2.15%	0.159%
32	11.027	1519	1520	1526	rVV	72931	91245	2.38%	0.176%
33	11.074	1526	1528	1532	rVV	162555	151900	3.96%	0.292%
34	11.121	1532	1536	1539	rVV3	55271	86853	2.26%	0.167%

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

35	11.163	1541	1543	1549	rVB	229294	226793	5.91%	0.436%
36	11.268	1557	1561	1563	rBV	1340241	1221457	31.81%	2.350%
37	11.292	1563	1565	1569	rVV	2466005	2101876	54.74%	4.043%
38	11.345	1569	1574	1577	rVB	705894	621200	16.18%	1.195%
39	11.380	1577	1580	1583	rBV2	90902	99520	2.59%	0.191%
40	11.498	1597	1600	1603	rBV	127894	129898	3.38%	0.250%
41	11.563	1608	1611	1614	rVV	111619	93516	2.44%	0.180%
42	11.598	1614	1617	1622	rVV2	159603	156078	4.06%	0.300%
43	11.680	1629	1631	1635	rVB	115625	114455	2.98%	0.220%
44	11.727	1636	1639	1641	rBV	62676	61663	1.61%	0.119%
45	11.768	1643	1646	1649	rVV	542886	514305	13.39%	0.989%
46	11.798	1649	1651	1655	rVV	722072	592518	15.43%	1.140%
47	11.845	1656	1659	1662	rVV	247718	208119	5.42%	0.400%
48	11.880	1662	1665	1668	rVV	800318	907055	23.62%	1.745%
49	11.904	1668	1669	1674	rVB	388672	237911	6.20%	0.458%
50	12.063	1693	1696	1699	rVV	305017	279995	7.29%	0.539%
51	12.092	1699	1701	1707	rVB	220130	223796	5.83%	0.431%
52	12.198	1716	1719	1720	rBV	70785	75557	1.97%	0.145%
53	12.215	1720	1722	1724	rVV	133249	122315	3.19%	0.235%
54	12.245	1724	1727	1729	rVV	181942	171054	4.45%	0.329%
55	12.268	1729	1731	1734	rVB2	135650	112314	2.93%	0.216%
56	12.321	1736	1740	1742	rBV	388367	392110	10.21%	0.754%
57	12.357	1742	1746	1748	rVV3	218604	279276	7.27%	0.537%
58	12.374	1748	1749	1752	rVB	106555	76866	2.00%	0.148%
59	12.410	1752	1755	1759	rBV	289385	329518	8.58%	0.634%
60	12.486	1763	1768	1771	rBV	3737991	3391113	88.32%	6.524%
61	12.545	1776	1778	1781	rVB2	112325	91450	2.38%	0.176%
62	12.574	1781	1783	1786	rBV	144115	123444	3.21%	0.237%
63	12.633	1790	1793	1797	rVB2	106737	106896	2.78%	0.206%
64	12.715	1800	1807	1810	rBV	3268620	3839755	100.00%	7.387%
65	12.762	1812	1815	1819	rVB2	156828	222870	5.80%	0.429%
66	12.827	1823	1826	1827	rBV2	222212	187801	4.89%	0.361%
67	12.851	1827	1830	1832	rVB	687551	612025	15.94%	1.177%
68	12.927	1838	1843	1847	rBV2	342054	571369	14.88%	1.099%
69	12.986	1850	1853	1854	rVV3	78126	77148	2.01%	0.148%
70	13.015	1854	1858	1859	rVV2	314444	361949	9.43%	0.696%
71	13.039	1859	1862	1865	rVB	611085	615271	16.02%	1.184%

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 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

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72	13.104	1870	1873	1876	rVV	381190	377415	9.83%	0.726%
73	13.139	1876	1879	1887	rVB3	478016	679459	17.70%	1.307%
74	13.198	1887	1889	1892	rBV3	85092	95836	2.50%	0.184%
75	13.233	1892	1895	1898	rVV	292237	256086	6.67%	0.493%
76	13.262	1898	1900	1904	rVV	267971	271335	7.07%	0.522%
77	13.551	1946	1949	1953	rVB	351976	356020	9.27%	0.685%
78	13.657	1963	1967	1969	rBV	454637	463112	12.06%	0.891%
79	13.680	1969	1971	1974	rVV	330713	337009	8.78%	0.648%
80	13.709	1974	1976	1981	rVV2	304463	454455	11.84%	0.874%
81	13.757	1981	1984	1988	rVB	402221	411567	10.72%	0.792%
82	13.827	1993	1996	1998	rBV	91112	102951	2.68%	0.198%
83	13.904	2004	2009	2012	rBV2	2310724	3244232	84.49%	6.241%
84	13.939	2012	2015	2018	rVB	1945834	1700486	44.29%	3.271%
85	14.015	2025	2028	2030	rVB	322512	250256	6.52%	0.481%
86	14.057	2032	2035	2040	rBV2	240785	308728	8.04%	0.594%
87	14.133	2044	2048	2051	rBV2	134786	216643	5.64%	0.417%
88	14.257	2066	2069	2071	rBV2	161214	144742	3.77%	0.278%
89	14.292	2073	2075	2078	rVB	429014	389334	10.14%	0.749%
90	14.327	2078	2081	2084	rBV	227282	193764	5.05%	0.373%
91	14.421	2094	2097	2098	rBV	274010	267656	6.97%	0.515%
92	14.780	2155	2158	2161	rBV2	168853	186117	4.85%	0.358%
93	14.939	2179	2185	2191	rBV	1796970	3551500	92.49%	6.832%
94	15.033	2198	2201	2205	rBV	345996	377146	9.82%	0.726%
95	15.221	2229	2233	2239	rVB	990477	1268907	33.05%	2.441%
96	15.286	2239	2244	2249	rBV	1530688	1867527	48.64%	3.593%
97	15.339	2250	2253	2256	rBV	683161	786768	20.49%	1.514%
98	15.368	2256	2258	2265	rVB2	345428	470659	12.26%	0.905%
99	16.750	2487	2493	2499	rVB	627233	1146063	29.85%	2.205%
100	17.180	2560	2566	2581	rVB	454625	1013107	26.38%	1.949%

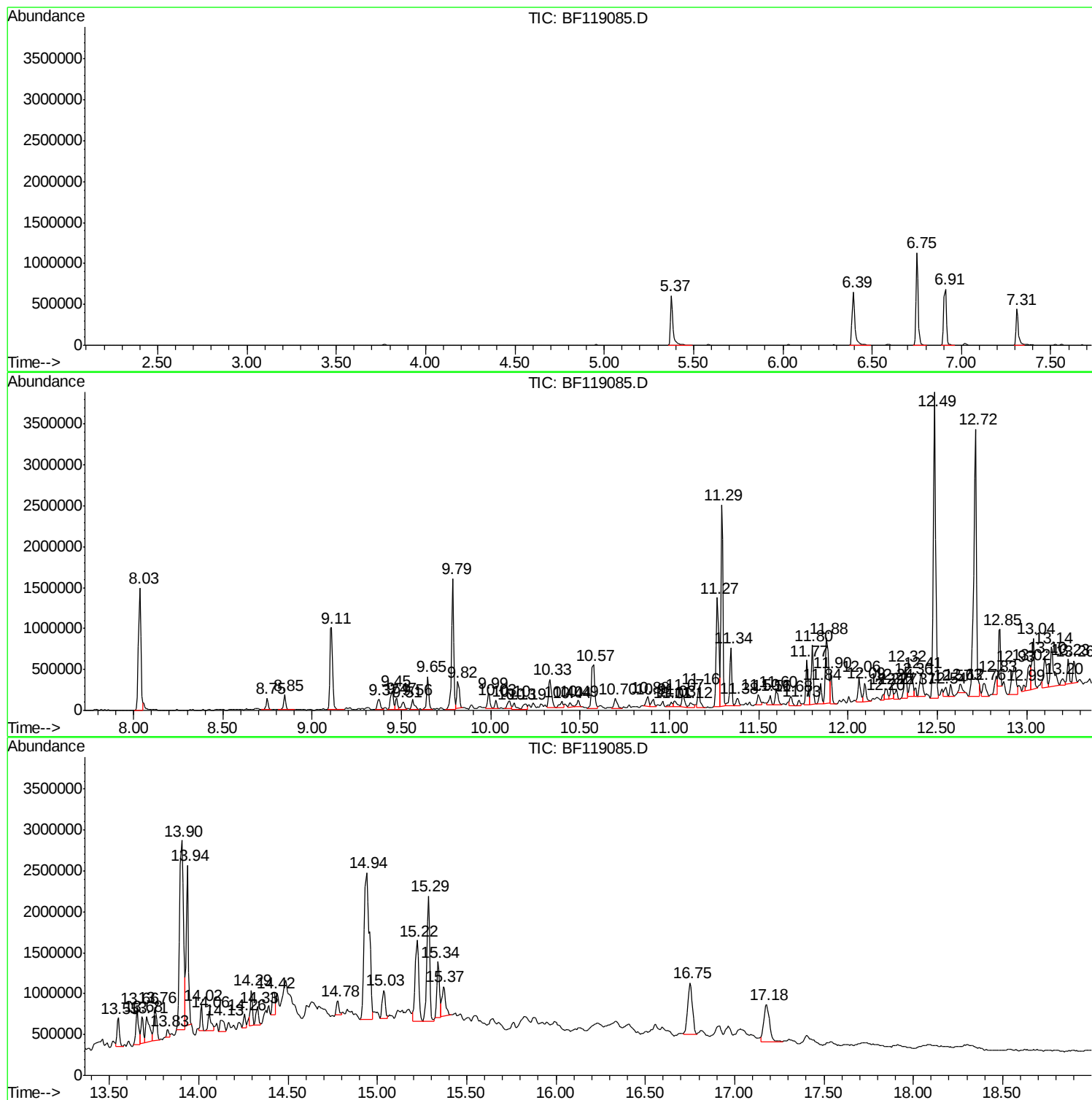
Sum of corrected areas: 51981963

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Instrument :
BNA_F
ClientSampled :

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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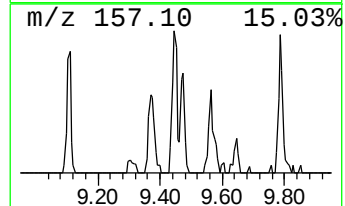
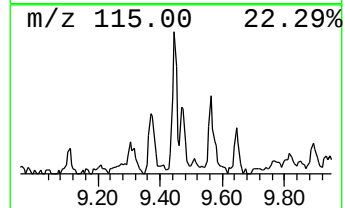
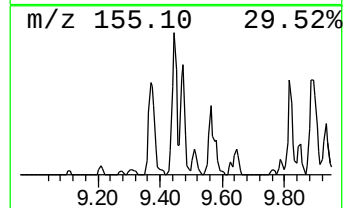
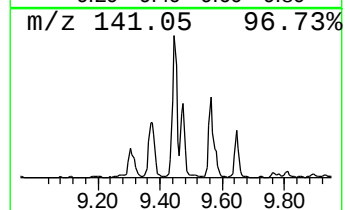
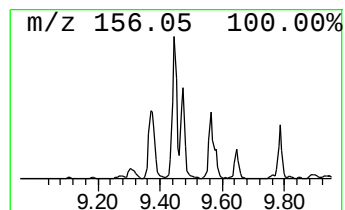
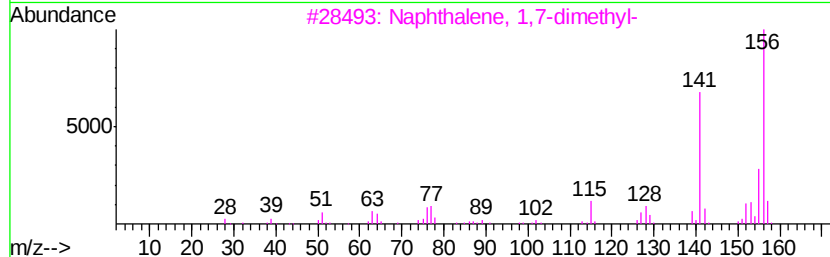
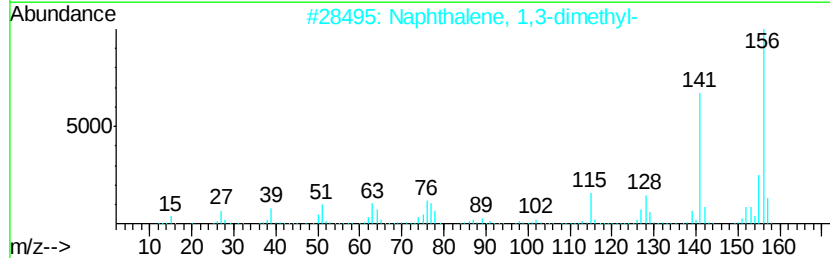
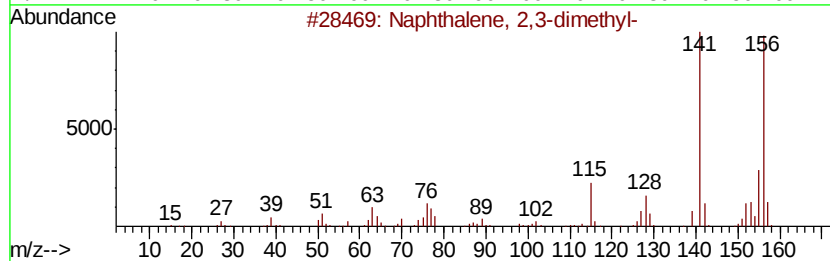
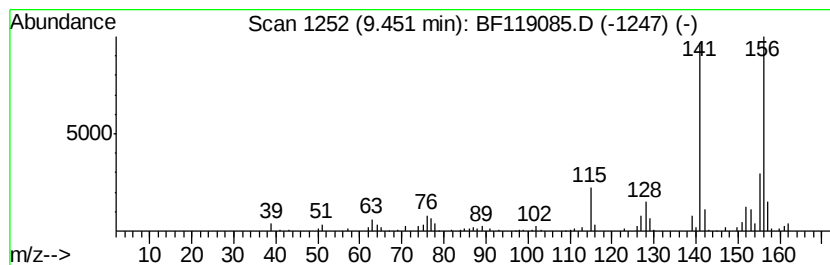
Instrument :
BNA_F
ClientSampleId :

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, 2,3-dimethyl- Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.45	3.32 ng	236697	Acenaphthene-d10	9.79
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 2,3-dimethyl-	156 C12H12	000581-40-8 98
2		Naphthalene, 1,3-dimethyl-	156 C12H12	000575-41-7 97
3		Naphthalene, 1,7-dimethyl-	156 C12H12	000575-37-1 97
4		Naphthalene, 1,6-dimethyl-	156 C12H12	000575-43-9 96
5		Naphthalene, 1,4-dimethyl-	156 C12H12	000571-58-4 96



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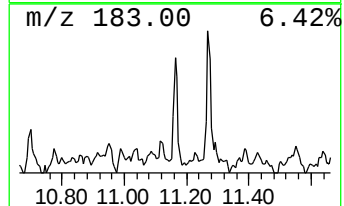
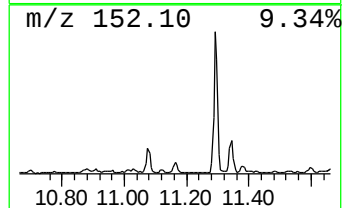
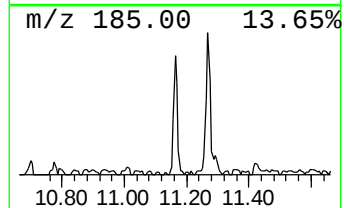
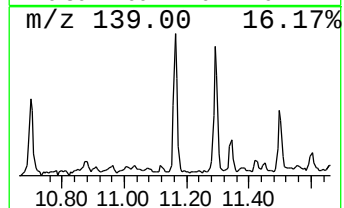
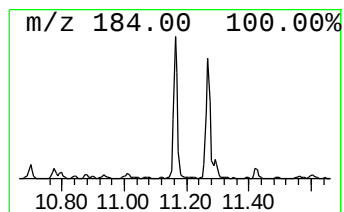
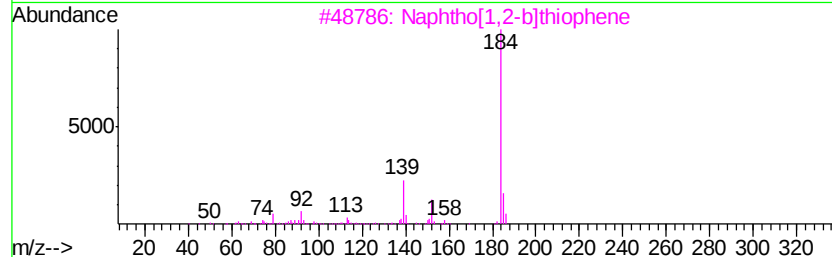
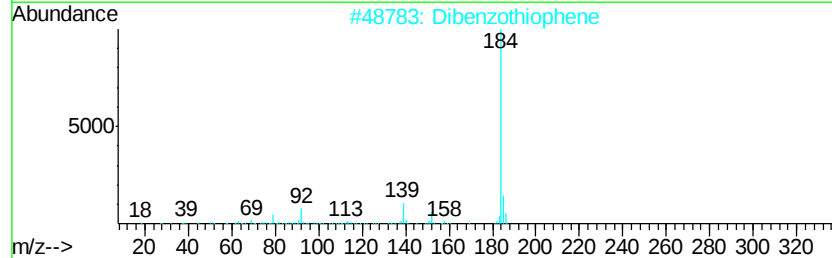
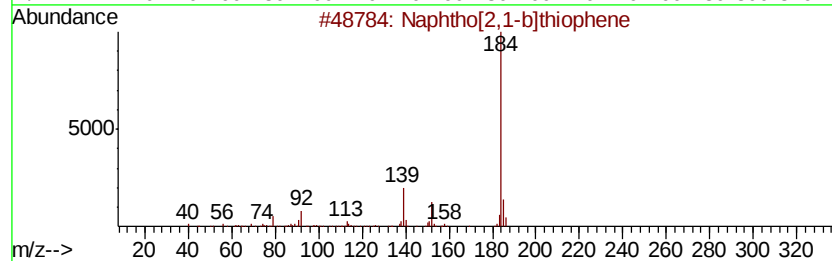
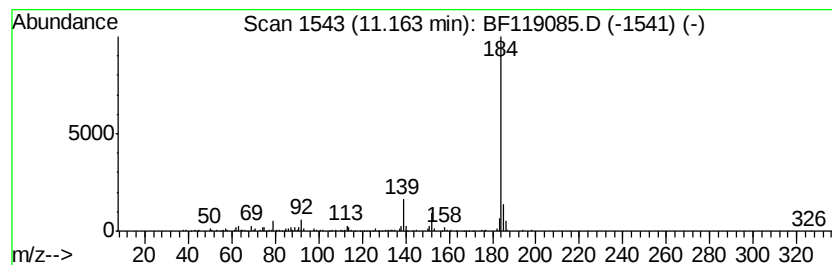
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 Naphtho[2,1-b]thiophene Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.16	3.71 ng	226793	Phenanthrene-d10	11.27

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphtho[2,1-b]thiophene	184	C12H8S	000233-02-3	95
2		Dibenzothiophene	184	C12H8S	000132-65-0	95
3		Naphtho[1,2-b]thiophene	184	C12H8S	000234-41-3	94
4		Naphtho[2,3-b]thiophene	184	C12H8S	000268-77-9	93
5		Azuleno(2,1-b)thiophene	184	C12H8S	000248-13-5	91



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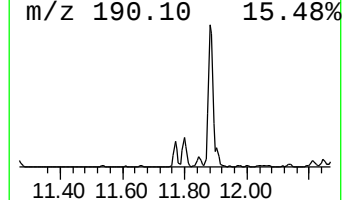
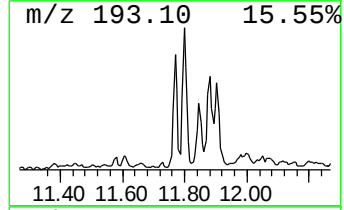
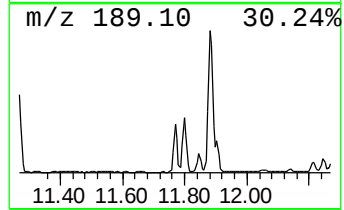
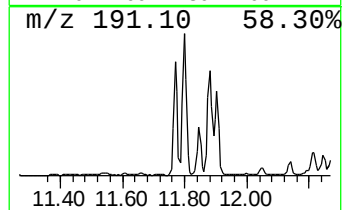
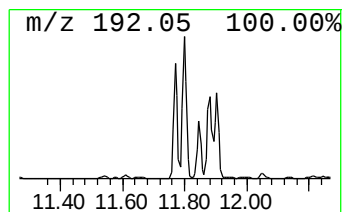
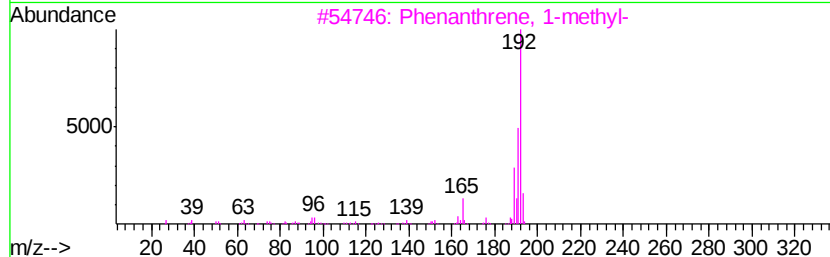
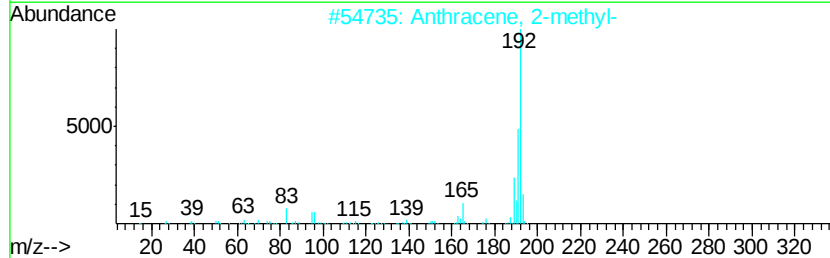
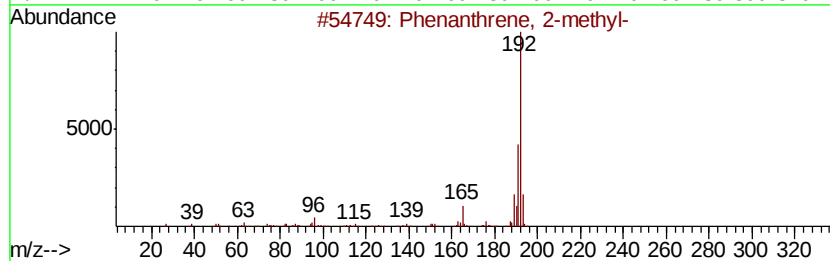
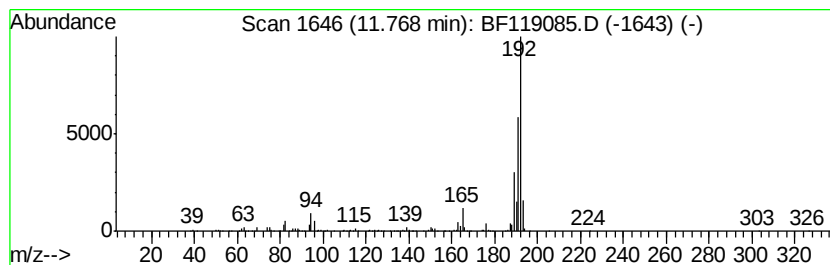
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Phenanthrene, 2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.77	8.42 ng	514305	Phenanthrene-d10	11.27

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	96
2		Anthracene, 2-methyl-	192	C15H12	000613-12-7	96
3		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	95
4		1H-Cyclopropa[1,1]phenanthrene, 1a, ...	192	C15H12	000949-41-7	94
5		Anthracene, 1-methyl-	192	C15H12	000610-48-0	94



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF022520\
Data File : BF119085.D
Acq On : 26 Feb 2020 1:10
Operator : CG/JU
Sample : L1367-05 5X
Misc :
ALS Vial : 23 Sample Multiplier: 1

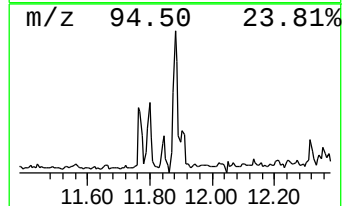
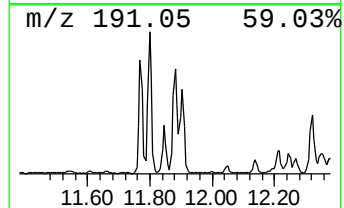
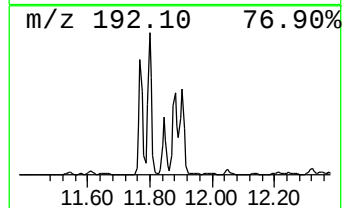
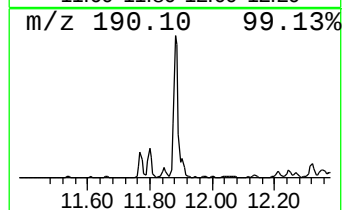
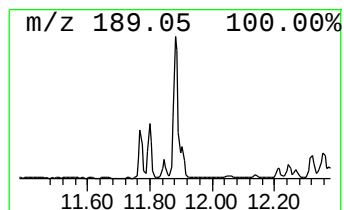
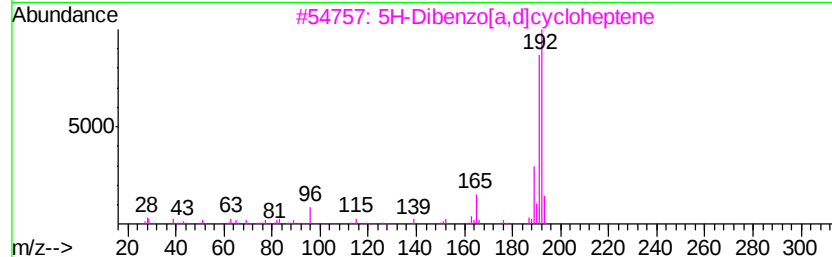
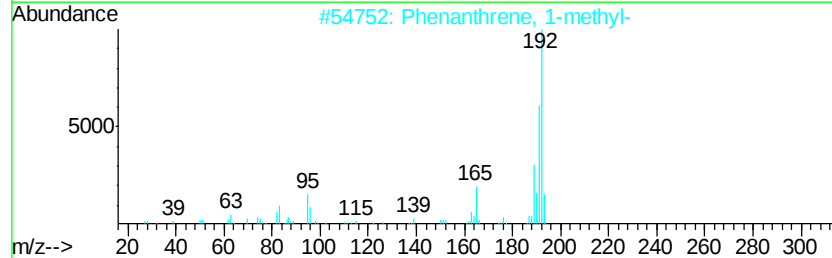
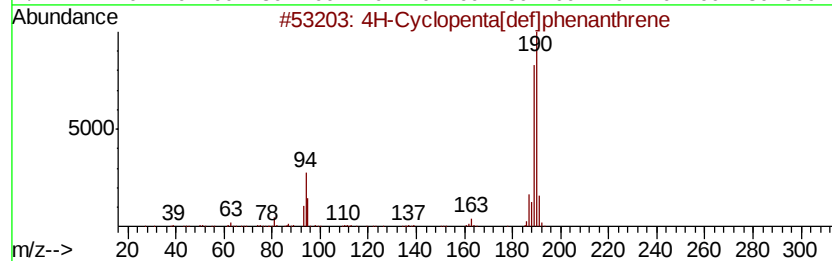
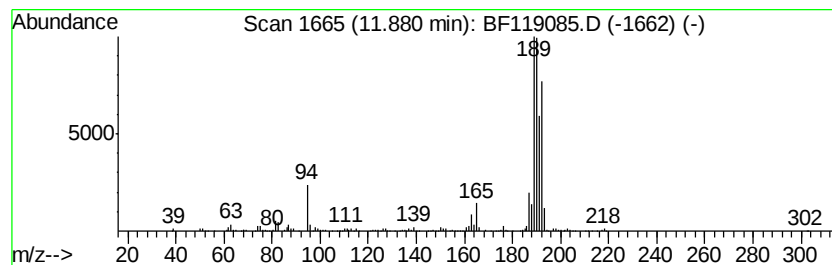
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 4H-Cyclopenta[def]phenanthrene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD		R.T.
11.88	14.85 ng	907055	Phenanthrene-d10		11.27
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	60
2	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	42
3	5H-Dibenzo[a,d]cycloheptene	192	C15H12	000256-81-5	42
4	Propyne, 1,3-diphenyl-	192	C15H12	004980-70-5	38
5	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	38



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF022520\
 Data File : BF119085.D
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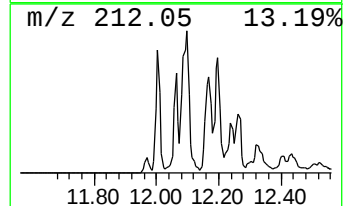
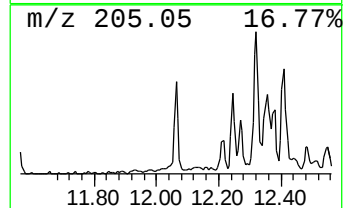
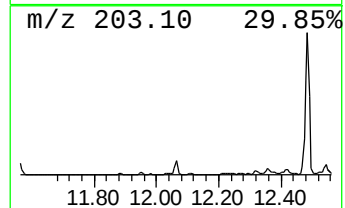
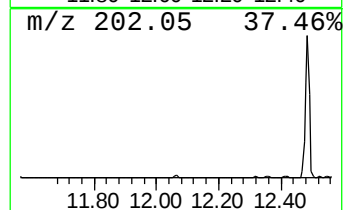
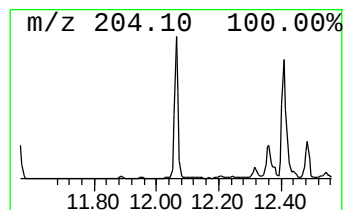
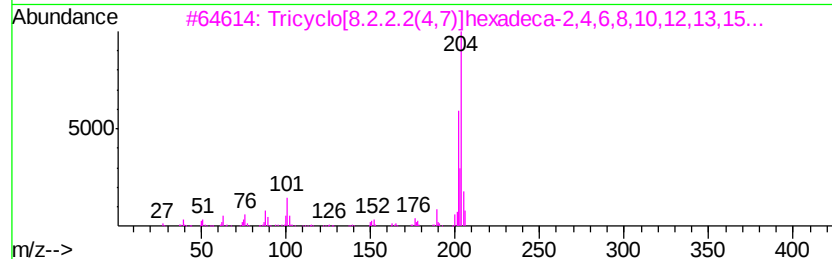
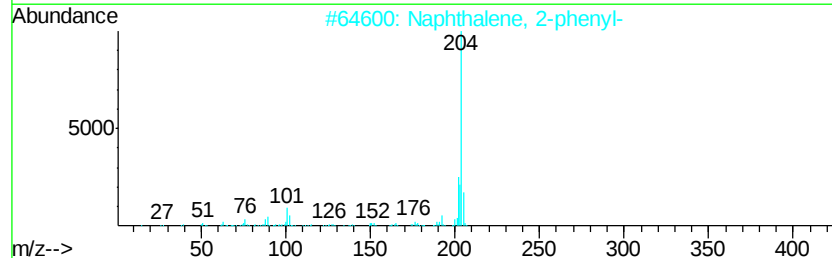
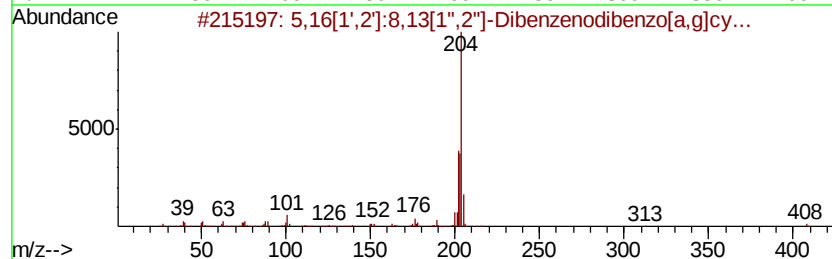
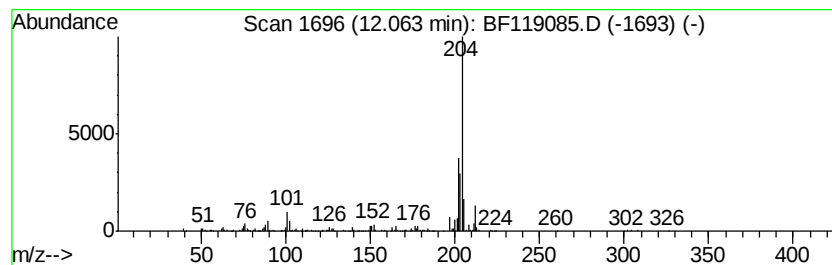
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 9 5,16[1',2']:8,13[1'',2'']-D... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.06	4.58 ng	279995	Phenanthrene-d10	11.27

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5,16[1',2']:8,13[1'',2'']-Dibenz...	408	C32H24	005672-97-9	80
2		Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	76
3		Tricyclo[8.2.2.2(4,7)]hexadeca-2...	204	C16H12	006572-60-7	68
4		9,10-Bis(bromomethyl)anthracene	362	C16H12Br2	034373-96-1	58
5		Levamisole	204	C11H12N2S	014769-73-4	52



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF022520\
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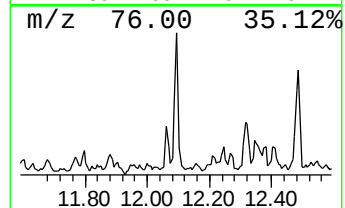
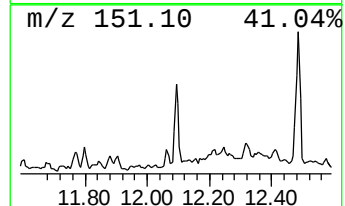
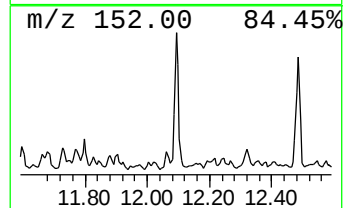
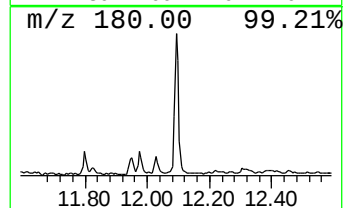
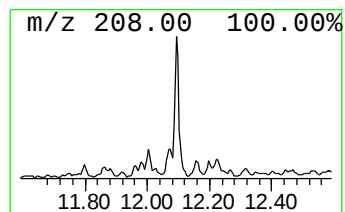
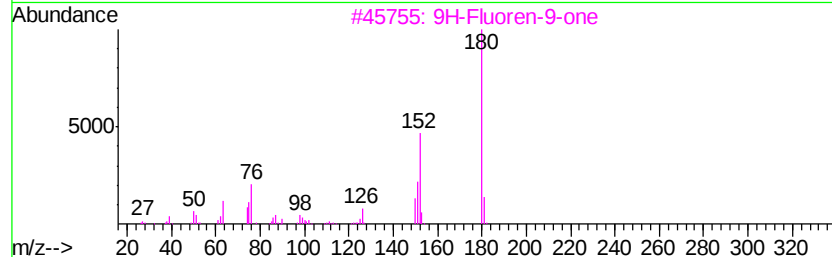
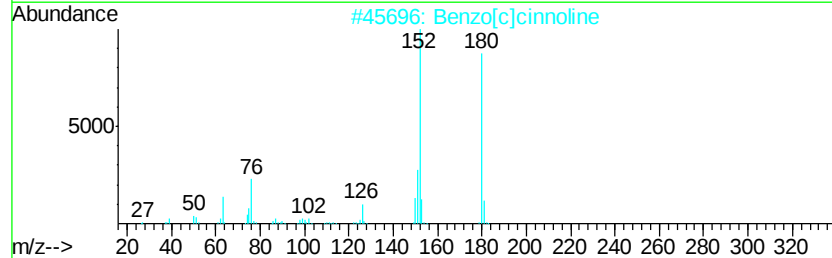
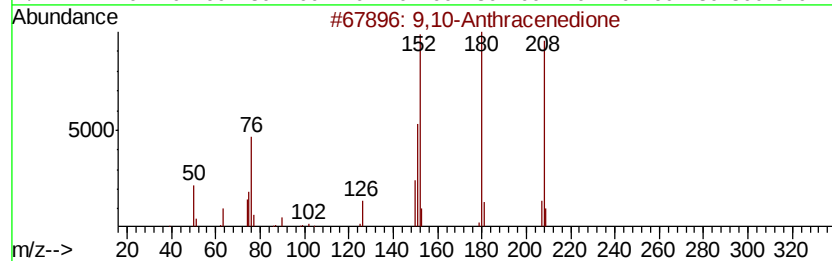
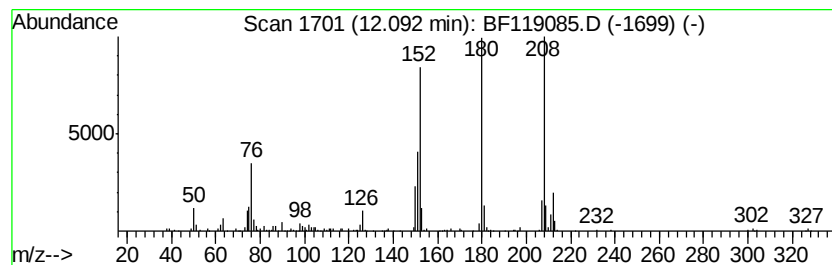
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 9,10-Anthracenedione Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.09	3.66 ng	223796	Phenanthrene-d10	11.27

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			9,10-Anthracenedione	208	C14H8O2	000084-65-1	99
2			Benzo[c]cinnoline	180	C12H8N2	000230-17-1	60
3			9H-Fluoren-9-one	180	C13H8O	000486-25-9	58
4			1H-Phenalen-1-one	180	C13H8O	000548-39-0	55
5			1,4-Anthracenedione	208	C14H8O2	000635-12-1	49



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF022520\
Data File : BF119085.D
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Sample : L1367-05 5X
Misc :
ALS Vial : 23 Sample Multiplier: 1

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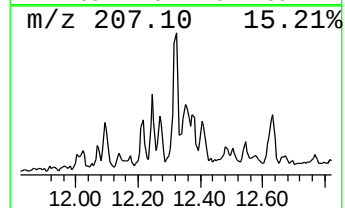
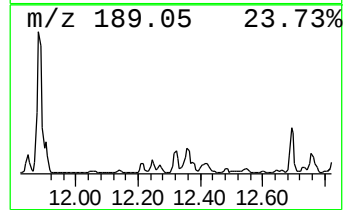
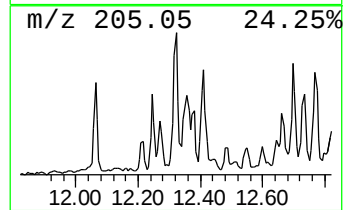
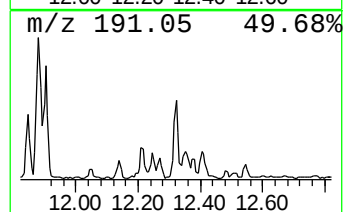
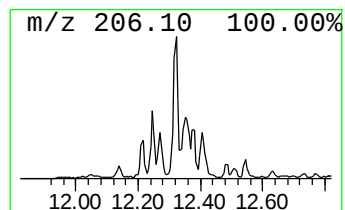
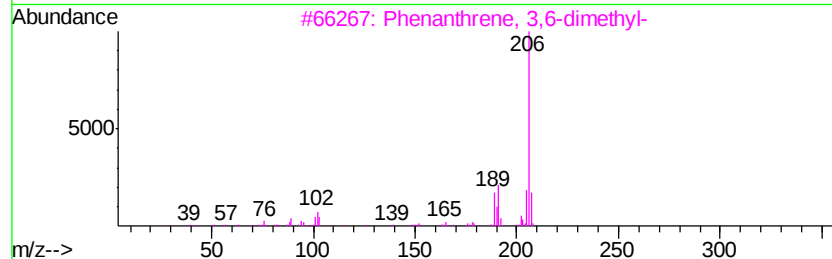
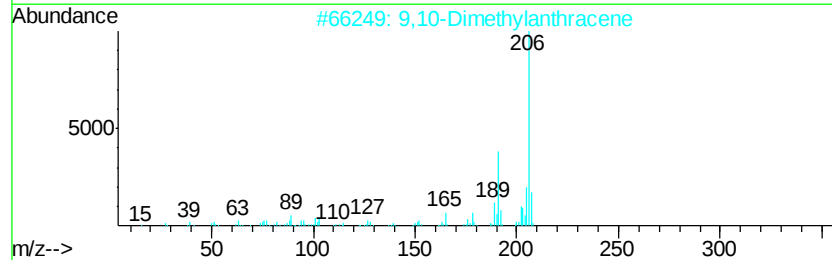
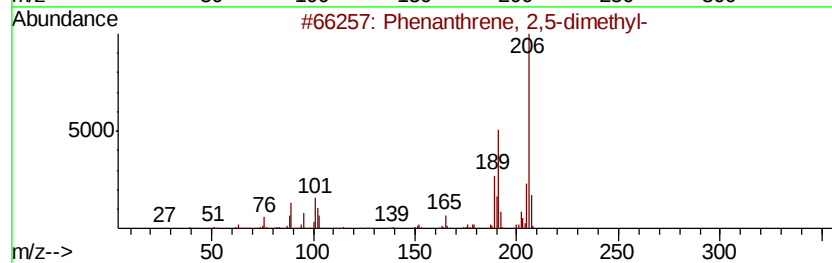
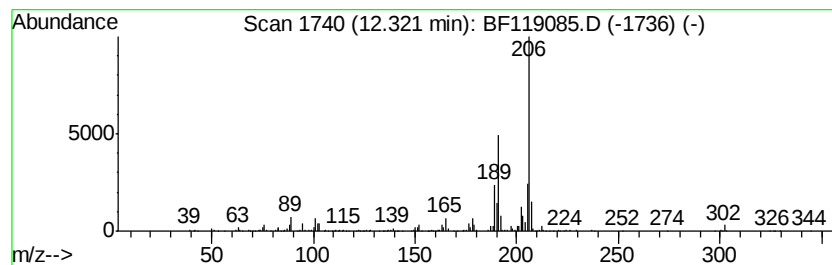
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Phenanthrene, 2,5-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.32	6.42 ng	392110	Phenanthrene-d10	11.27

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	96
2			9,10-Dimethylantracene	206	C16H14	000781-43-1	95
3			Phenanthrene, 3,6-dimethyl-	206	C16H14	001576-67-6	87
4			Phenanthrene, 2,3-dimethyl-	206	C16H14	003674-65-5	87
5			1,3-Diphenyl-3-methylcyclopropene	206	C16H14	086544-79-8	81



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF022520\
Data File : BF119085.D
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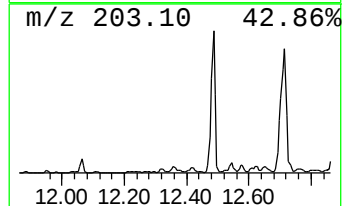
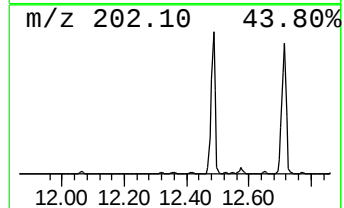
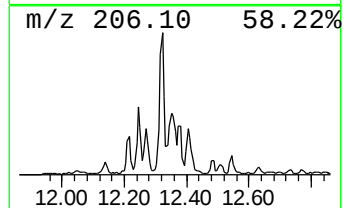
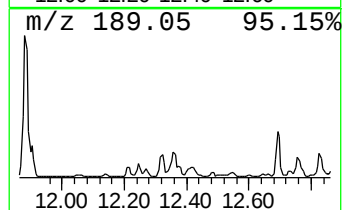
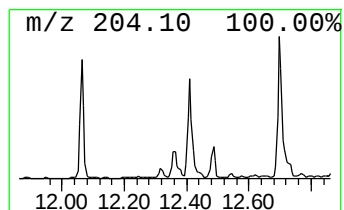
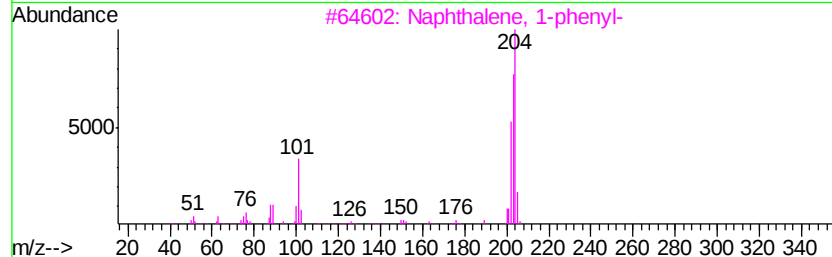
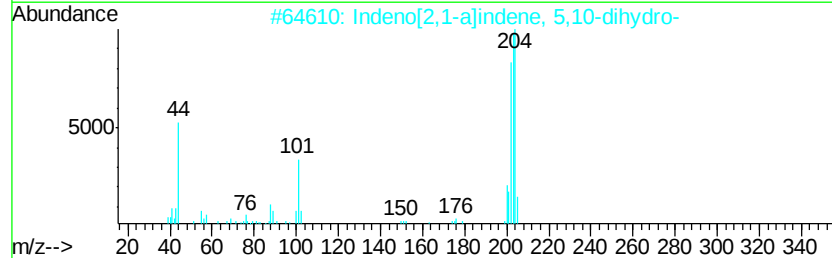
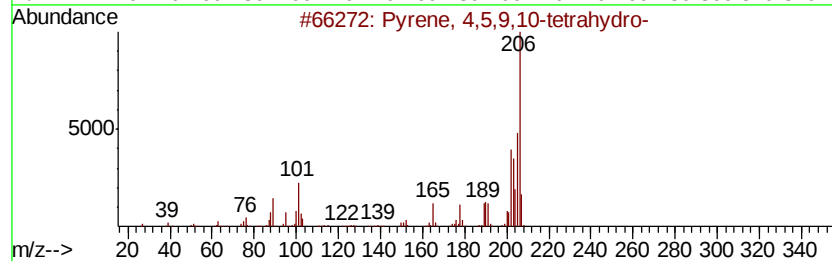
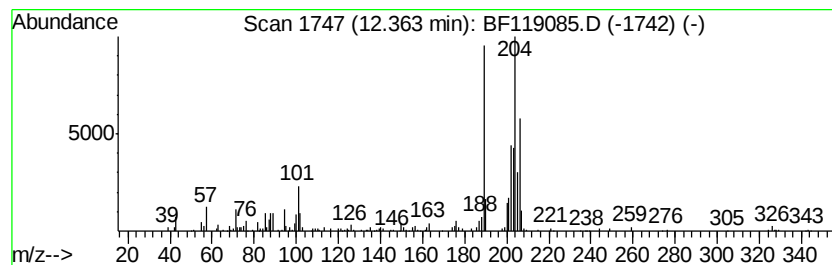
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 unknown12.36 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.36	4.57 ng	279276	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		Indeno[2,1-a]indene, 5,10-dihydro-	204	C16H12	006543-29-9	42
3		Naphthalene, 1-phenyl-	204	C16H12	000605-02-7	38
4		9,10-Bis(bromomethyl)anthracene	362	C16H12Br2	034373-96-1	35
5		1,2,3,3a-Tetrahydro-10-phenylben...	260	C18H16N2	028749-06-6	35



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF022520\
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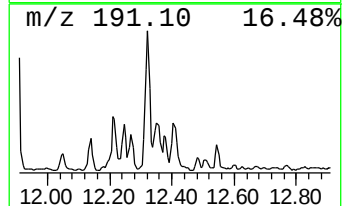
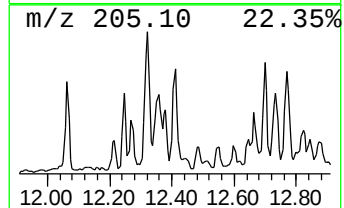
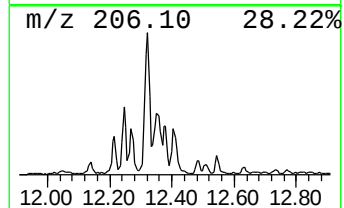
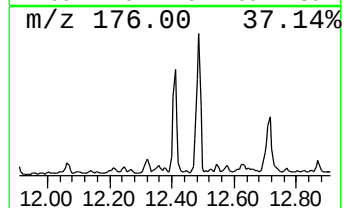
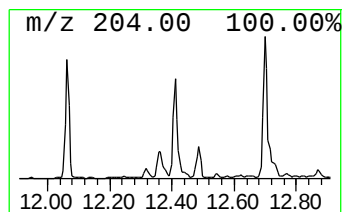
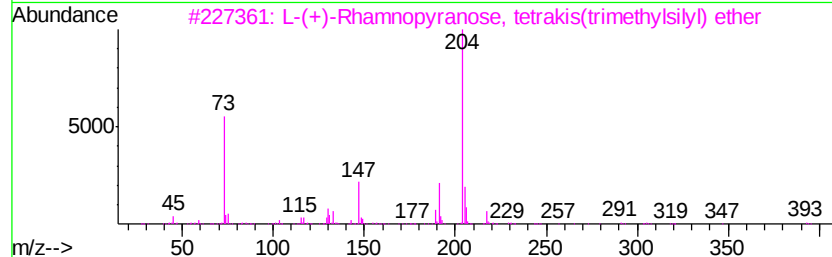
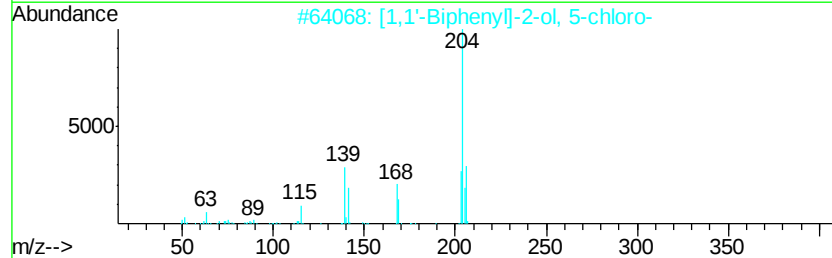
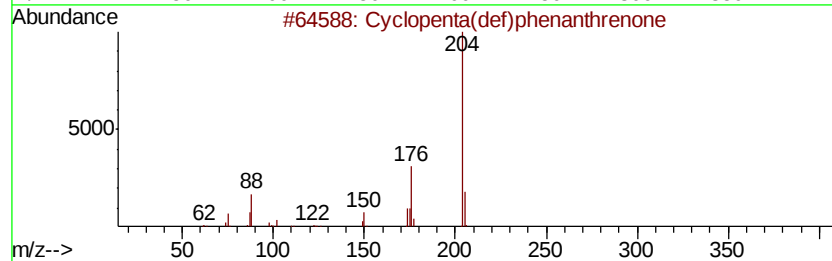
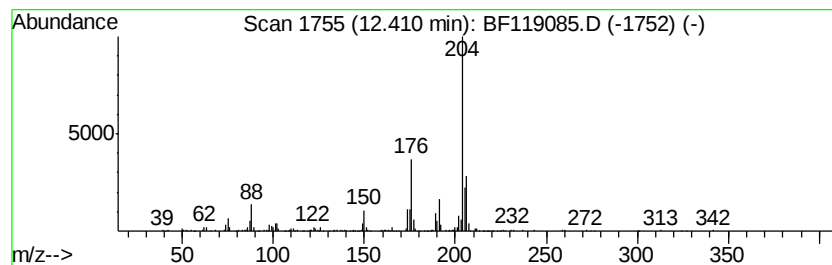
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Cyclopenta(def)phenanthrenone Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.41	5.40 ng	329518	Phenanthrene-d10	11.27

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopenta(def)phenanthrenone	204	C15H8O	005737-13-3	93
2		[1,1'-Biphenyl]-2-ol, 5-chloro-	204	C12H9ClO	000607-12-5	47
3		L-(+)-Rhamnopyranose, tetrakis(t...	452	C18H44O5Si4	1000380-12-9	47
4		1,2,4,8-Tetramethylbicyclo[6.3.0...	204	C15H24	137235-51-9	46
5		Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	43



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF022520\
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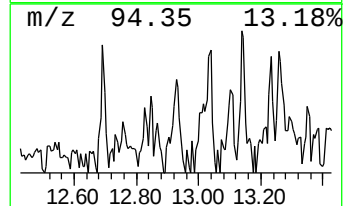
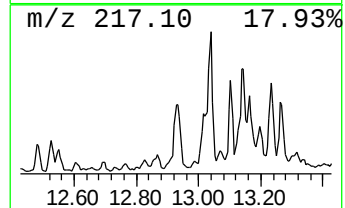
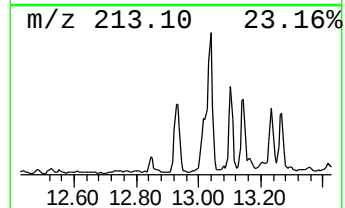
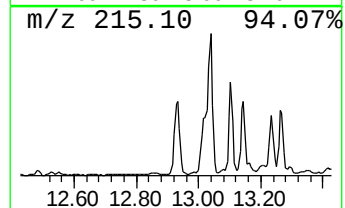
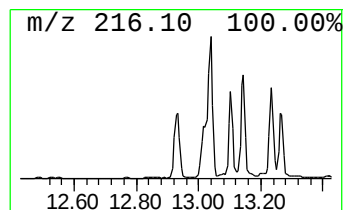
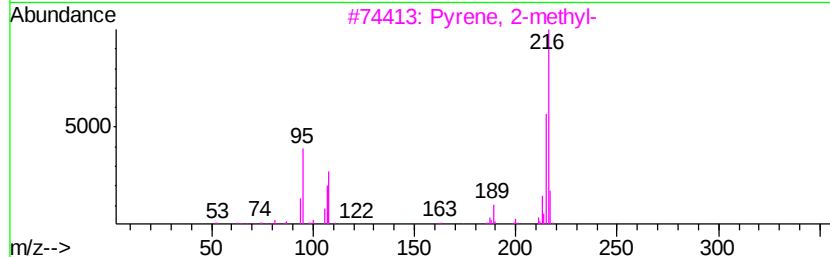
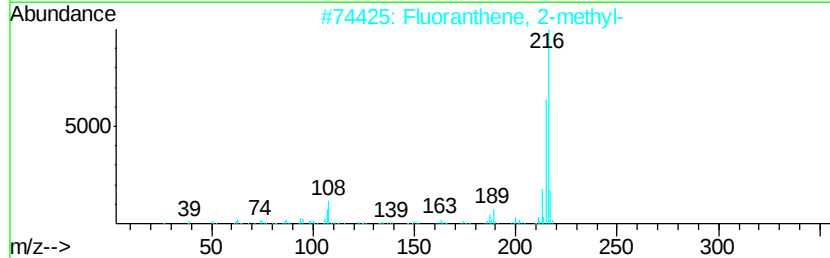
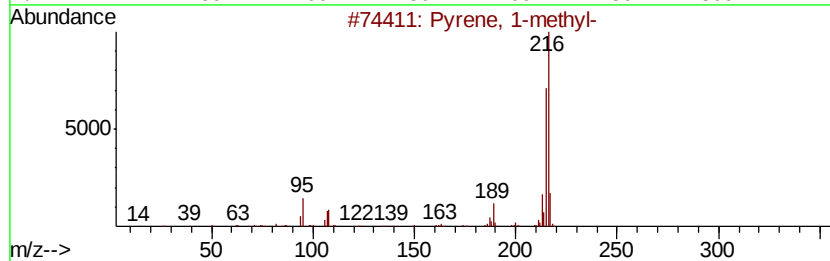
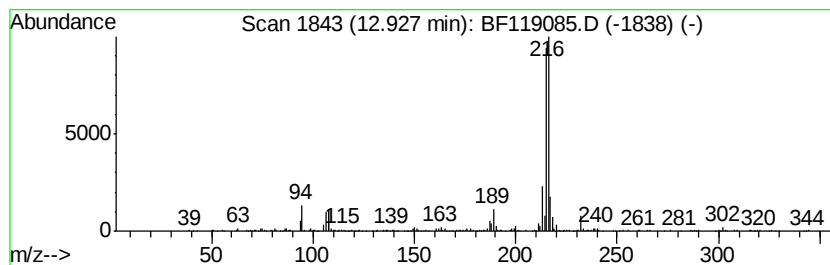
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ClientSampleId :

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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 14 Pyrene, 1-methyl- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.93	3.52 ng	571369	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 92
3		Pyrene, 2-methyl-	216 C17H12	003442-78-2 91
4		11H-Benzo[b]fluorene	216 C17H12	000243-17-4 90
5		7H-Benzo[c]fluorene	216 C17H12	000205-12-9 83



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF022520\
Data File : BF119085.D
Acq On : 26 Feb 2020 1:10
Operator : CG/JU
Sample : L1367-05 5X
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
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ClientSampleId :

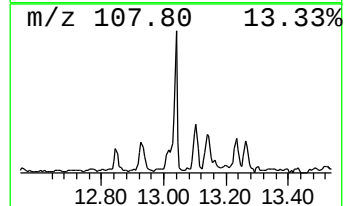
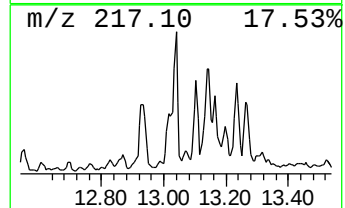
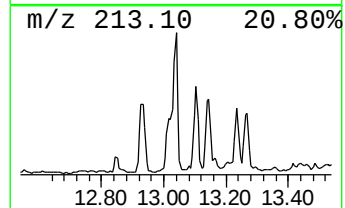
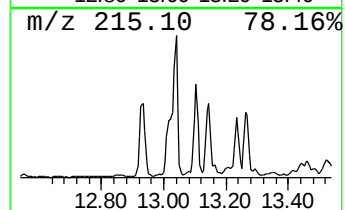
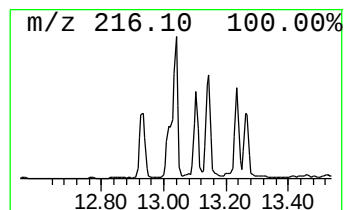
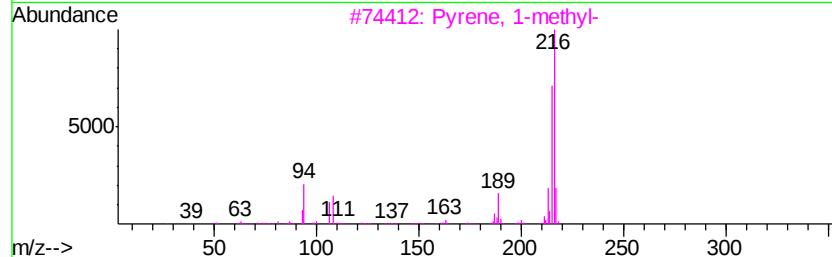
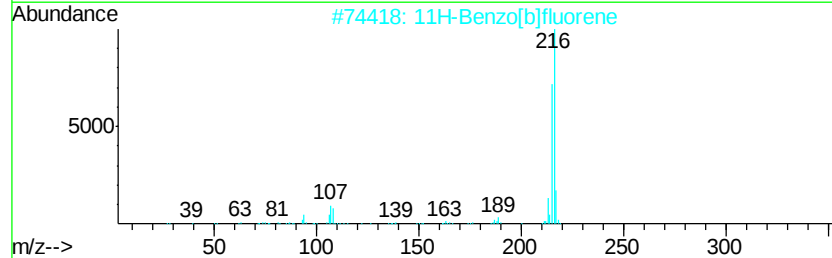
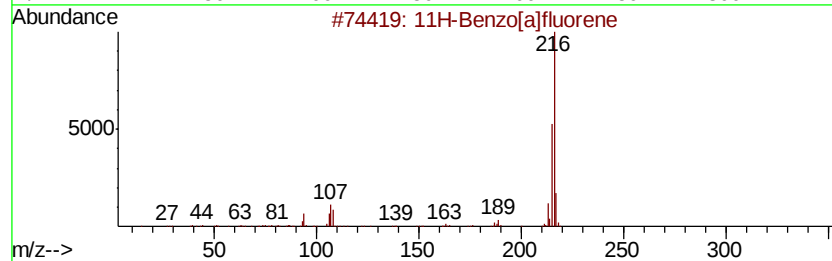
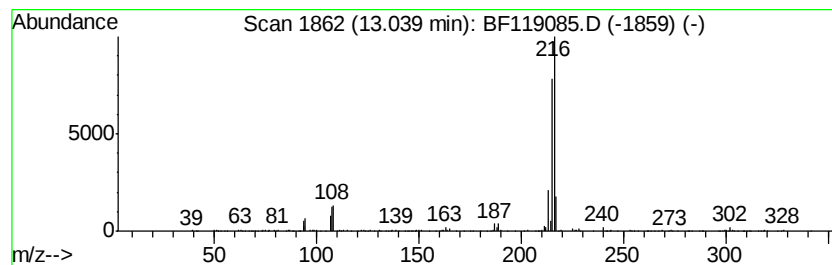
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 11H-Benzo[a]fluorene Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.04	3.79 ng	615271	Chrysene-d12	13.91

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	94
2		11H-Benzo[b]fluorene	216	C17H12	000243-17-4	93
3		Pvrene, 1-methyl-	216	C17H12	002381-21-7	74
4		6H-Thiazolo[5,4-e]indole, 2,7,8-...	216	C12H12N2S	1000338-32-0	59
5		Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	53



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF022520\
Data File : BF119085.D
Acq On : 26 Feb 2020 1:10
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Sample : L1367-05 5X
Misc :
ALS Vial : 23 Sample Multiplier: 1

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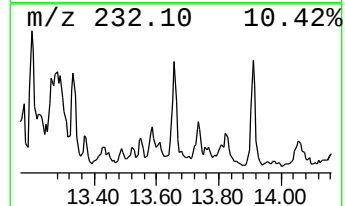
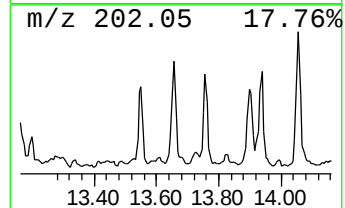
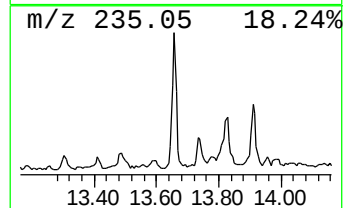
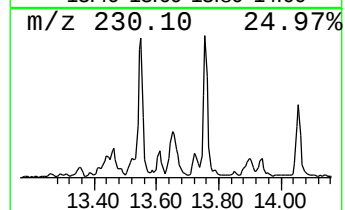
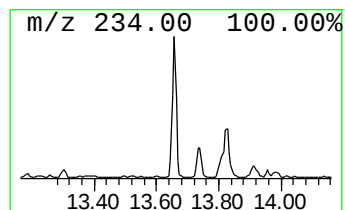
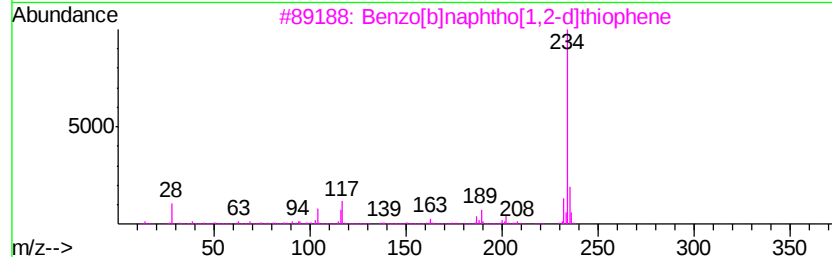
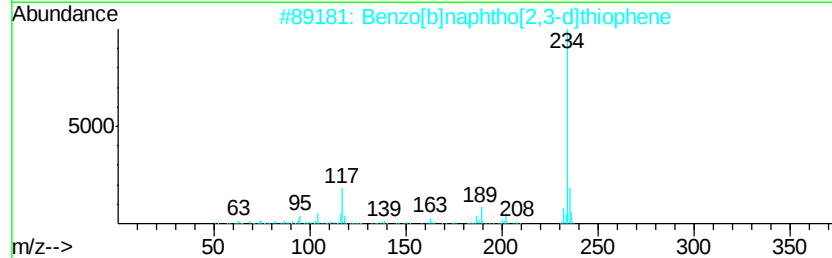
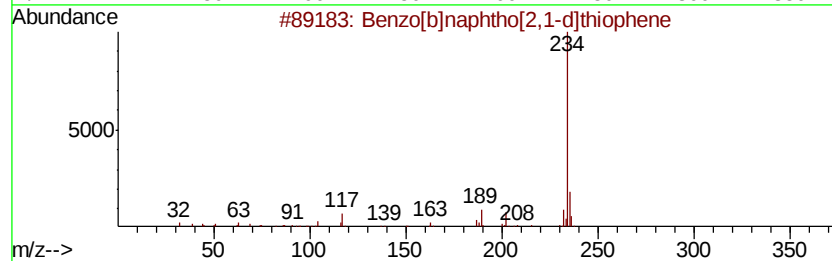
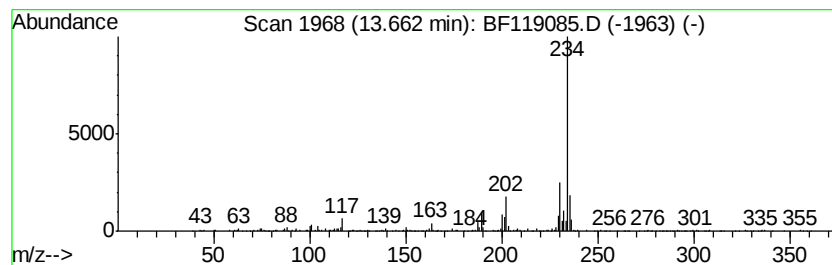
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 Benzo[b]naphtho[2,1-d]thiop... Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.66	2.85 ng	463112	Chrysene-d12	13.91

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzo[b]naphtho[2,1-d]thiophene	234	C16H10S	000239-35-0	93
2		Benzo[b]naphtho[2,3-d]thiophene	234	C16H10S	000243-46-9	62
3		Benzo[b]naphtho[1,2-d]thiophene	234	C16H10S	000205-43-6	58
4		2-Amino-6-t-butyl-3-cvano-4,5,6,...	234	C13H18N2S	042159-76-2	46
5		Quinoxalino[2,3-b]quinoxaline, 5...	234	C14H10N4	000531-46-4	43



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF022520\
Data File : BF119085.D
Acq On : 26 Feb 2020 1:10
Operator : CG/JU
Sample : L1367-05 5X
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
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ClientSampleId :

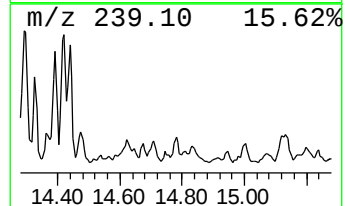
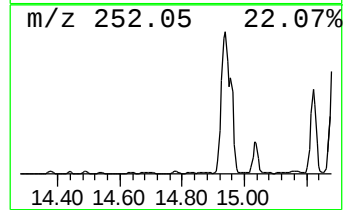
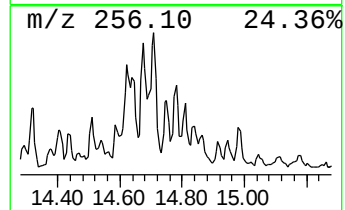
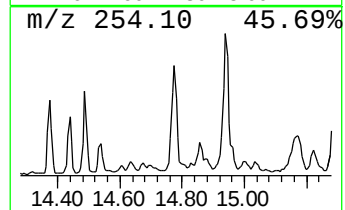
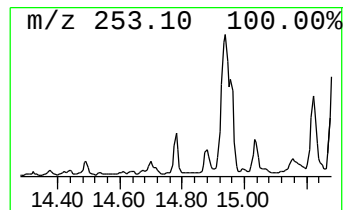
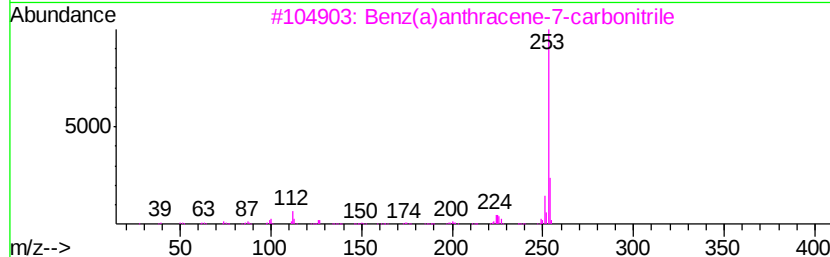
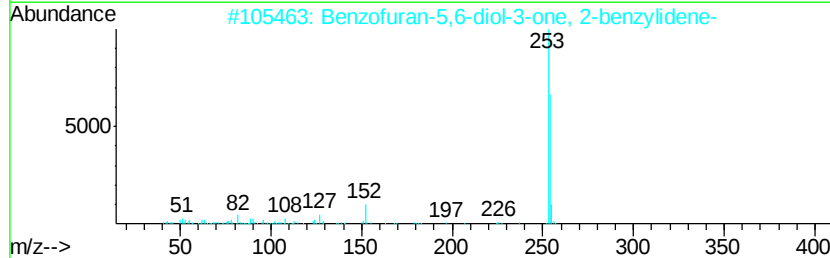
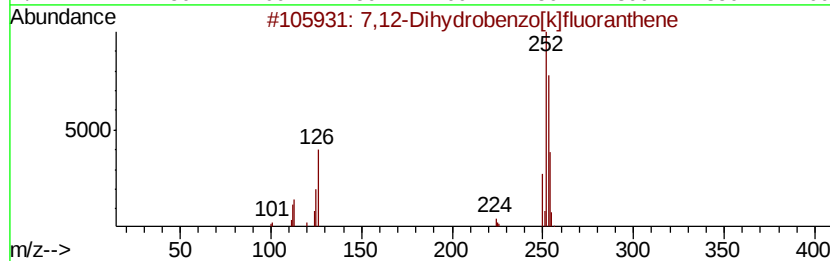
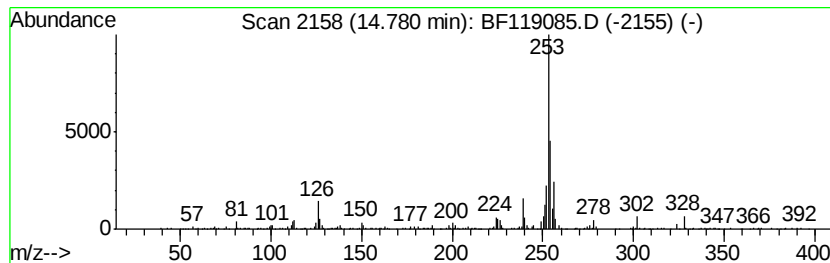
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 7,12-Dihydrobenzo[k]fluoran... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.78	4.73 ng	186117	Perylene-d12	15.34

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			7,12-Dihydrobenzo[k]fluoranthene	254	C20H14	1000080-17-7	62
2			Benzofuran-5,6-diol-3-one, 2-ben...	254	C15H10O4	1000128-65-2	50
3			Benz(a)anthracene-7-carbonitrile	253	C19H11N	007476-08-6	43
4			10-Oxo-5,5-dimethyl-5-sila-5,10-...	254	C14H14N2OSi	089052-45-9	42
5			4,6'-BiazulenyI	254	C20H14	094154-49-1	38



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF022520\
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Operator : CG/JU
Sample : L1367-05 5X
Misc :
ALS Vial : 23 Sample Multiplier: 1

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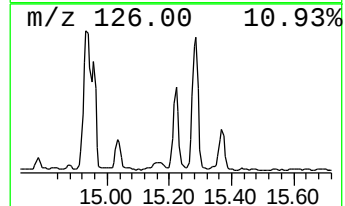
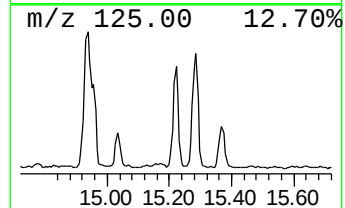
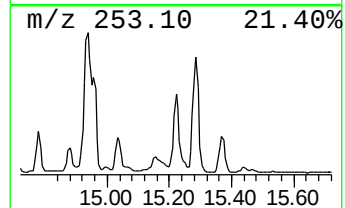
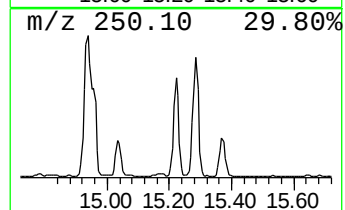
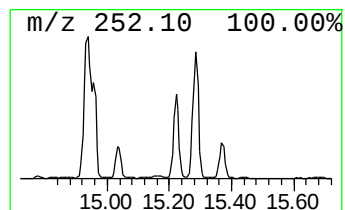
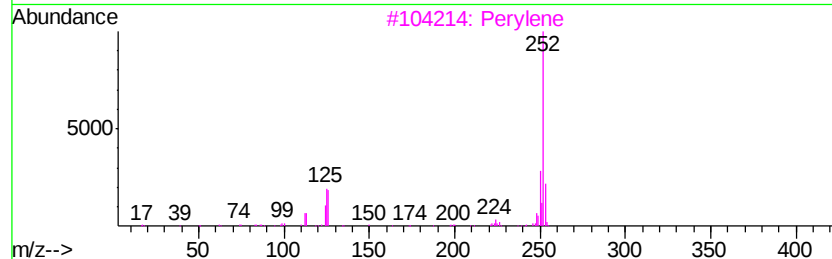
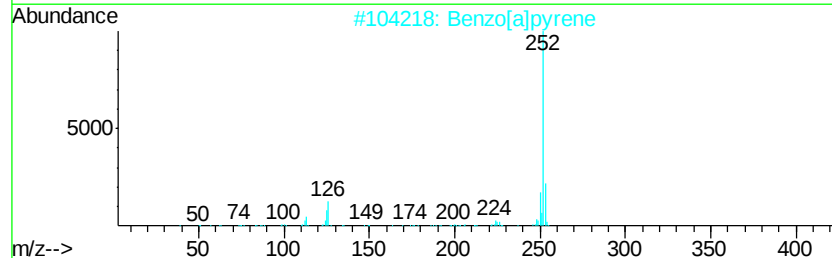
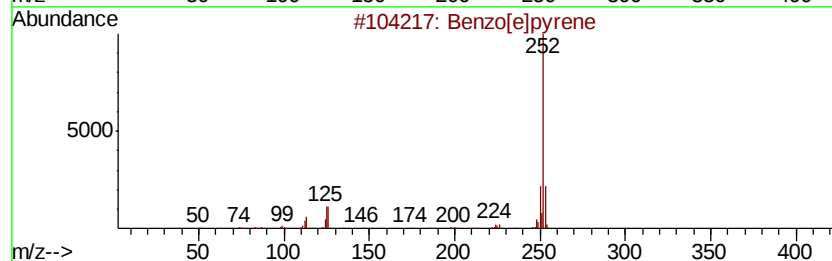
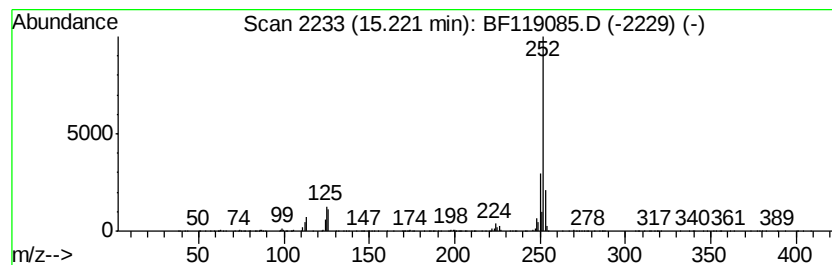
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 20 Benzo[e]pyrene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.22	32.26 ng	1268910	Perylene-d12	15.34

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



Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF022520\
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 Operator : CG/JU
 Sample : L1367-05 5X
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :

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, 2,3-...	9.45	3.3	ng	236697	3	9.79	1425550	20.0
Naphtho[2,1-b]thi...	11.16	3.7	ng	226793	4	11.27	1221460	20.0
Phenanthrene, 2-m...	11.77	8.4	ng	514305	4	11.27	1221460	20.0
4H-Cyclopenta[def...	11.88	14.9	ng	907055	4	11.27	1221460	20.0
5,16[1',2']:8,13[...	12.06	4.6	ng	279995	4	11.27	1221460	20.0
9,10-Anthracenedione	12.09	3.7	ng	223796	4	11.27	1221460	20.0
Phenanthrene, 2,5...	12.32	6.4	ng	392110	4	11.27	1221460	20.0
unknown12.36	12.36	4.6	ng	279276	4	11.27	1221460	20.0
Cyclopenta(def)ph...	12.41	5.4	ng	329518	4	11.27	1221460	20.0
Pyrene, 1-methyl-	12.93	3.5	ng	571369	5	13.91	3244230	20.0
11H-Benzo[a]fluorene	13.04	3.8	ng	615271	5	13.91	3244230	20.0
Benzo[b]naphtho[2...	13.66	2.9	ng	463112	5	13.91	3244230	20.0
7,12-Dihydrobenzo...	14.78	4.7	ng	186117	6	15.34	786768	20.0
Benzo[e]pyrene	15.22	32.3	ng	1268910	6	15.34	786768	20.0