

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF122217\
 Data File : BF101658.D
 Acq On : 22 Dec 2017 17:26
 Operator : SJ/JU
 Sample : I6683-09 2X
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EO-02-122017-B

Integration Parameters: rteint.p
 Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0
 Filtering: 5
 Min Area: 3 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF121417.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.998	493	495	507	rBV	75326	132466	1.03%	0.150%
2	5.404	561	564	583	rBV	1249045	1751274	13.66%	1.987%
3	6.428	734	738	757	rBV	1345684	1870320	14.59%	2.122%
4	6.557	757	760	766	rVV	1866857	1724290	13.45%	1.956%
5	6.787	796	799	814	rBV	969604	920636	7.18%	1.045%
6	6.939	822	825	840	rVB	1419157	1352779	10.55%	1.535%
7	7.357	892	896	913	rBV	939757	1197441	9.34%	1.359%
8	8.069	1014	1017	1019	rVV	1323626	1139606	8.89%	1.293%
9	8.092	1019	1021	1028	rVB	409696	446563	3.48%	0.507%
10	8.645	1112	1115	1119	rVB	195074	155499	1.21%	0.176%
11	8.781	1135	1138	1145	rBV	529028	479888	3.74%	0.545%
12	8.881	1150	1155	1161	rVB	599178	565508	4.41%	0.642%
13	9.139	1195	1199	1203	rBV	2113116	1761340	13.74%	1.999%
14	9.210	1208	1211	1214	rBV	402731	325034	2.53%	0.369%
15	9.410	1240	1245	1249	rBV	363284	479357	3.74%	0.544%
16	9.480	1253	1257	1260	rVV	494331	566048	4.41%	0.642%
17	9.510	1260	1262	1264	rVV	318317	334116	2.61%	0.379%
18	9.533	1264	1266	1273	rVB3	210151	365811	2.85%	0.415%
19	9.598	1273	1277	1282	rBV2	261676	314288	2.45%	0.357%
20	9.686	1288	1292	1295	rBV2	388864	367133	2.86%	0.417%
21	9.739	1299	1301	1305	rVB	542887	410883	3.20%	0.466%
22	9.822	1312	1315	1318	rVB2	1221150	1013855	7.91%	1.150%
23	9.857	1318	1321	1324	rVB	662738	581971	4.54%	0.660%
24	9.928	1329	1333	1337	rBV3	126387	192030	1.50%	0.218%
25	9.975	1337	1341	1343	rBV3	116494	161956	1.26%	0.184%
26	10.028	1347	1350	1353	rVB2	557174	528397	4.12%	0.600%
27	10.063	1353	1356	1360	rVB3	275390	266087	2.08%	0.302%
28	10.139	1365	1369	1372	rBV	228661	275989	2.15%	0.313%
29	10.169	1372	1374	1380	rVB2	183865	181320	1.41%	0.206%
30	10.239	1380	1386	1390	rBV2	640495	757644	5.91%	0.860%
31	10.375	1405	1409	1415	rVB	1009807	1009164	7.87%	1.145%
32	10.457	1415	1423	1425	rBV3	347750	545992	4.26%	0.620%
33	10.527	1433	1435	1439	rVB	195884	202101	1.58%	0.229%
34	10.616	1446	1450	1456	rBV2	987398	1089781	8.50%	1.237%

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Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF121417.M
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35	10.710	1463	1466	1473	rBV2	528591	691187	5.39%	0.784%
36	10.922	1496	1502	1505	rBV2	340818	467437	3.65%	0.530%
37	10.951	1505	1507	1509	rVV	222903	189394	1.48%	0.215%
38	11.004	1513	1516	1518	rVB2	224991	201233	1.57%	0.228%
39	11.051	1518	1524	1526	rBV3	137281	255165	1.99%	0.290%
40	11.157	1540	1542	1545	rBV	486758	460781	3.59%	0.523%
41	11.186	1545	1547	1549	rVV	168529	166495	1.30%	0.189%
42	11.210	1549	1551	1556	rVB	444785	384692	3.00%	0.436%
43	11.316	1565	1569	1571	rBV	1108134	1120973	8.74%	1.272%
44	11.345	1571	1574	1577	rVB	3195018	3010934	23.48%	3.416%
45	11.392	1577	1582	1586	rBV	1271170	1101257	8.59%	1.250%
46	11.557	1605	1610	1613	rBV2	267789	364232	2.84%	0.413%
47	11.586	1613	1615	1617	rBV	310771	225308	1.76%	0.256%
48	11.645	1623	1625	1629	rVB	205520	185924	1.45%	0.211%
49	11.698	1629	1634	1637	rVV	4899175	5205439	40.60%	5.906%
50	11.727	1637	1639	1644	rVB	110407	138031	1.08%	0.157%
51	11.816	1651	1654	1657	rBV	790452	758129	5.91%	0.860%
52	11.845	1657	1659	1663	rVB	971087	810749	6.32%	0.920%
53	11.892	1664	1667	1670	rBV	392783	380143	2.96%	0.431%
54	11.927	1670	1673	1676	rBV2	1177378	1325221	10.34%	1.504%
55	11.951	1676	1677	1681	rVB	529212	347639	2.71%	0.394%
56	11.992	1681	1684	1686	rBV	294664	248840	1.94%	0.282%
57	12.110	1699	1704	1708	rBV2	457521	474656	3.70%	0.539%
58	12.145	1708	1710	1715	rVB3	110970	133709	1.04%	0.152%
59	12.292	1732	1735	1737	rVB	215012	150062	1.17%	0.170%
60	12.316	1737	1739	1742	rVB2	174116	155147	1.21%	0.176%
61	12.398	1744	1753	1755	rBV2	5410823	12822117	100.00%	14.549%
62	12.486	1764	1768	1771	rVB	3034938	2762391	21.54%	3.134%
63	12.539	1772	1777	1780	rBV	3380652	3635706	28.35%	4.125%
64	12.686	1800	1802	1805	rBV	274879	276582	2.16%	0.314%
65	12.774	1810	1817	1822	rVV2	2804528	4006524	31.25%	4.546%
66	12.816	1822	1824	1829	rVB	260226	260407	2.03%	0.295%
67	12.898	1834	1838	1841	rBV2	1441741	1254615	9.78%	1.424%
68	12.921	1841	1842	1848	rVB2	101162	134445	1.05%	0.153%
69	12.986	1848	1853	1858	rBV4	354397	641821	5.01%	0.728%
70	13.092	1869	1871	1876	rVV	799844	867621	6.77%	0.984%
71	13.157	1880	1882	1885	rVB	708668	619841	4.83%	0.703%

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Sampling : 1 Min Area: 3 % of largest Peak
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Stop Thrs : 0 Peak Location: TOP

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72	13.204	1885	1890	1896	rVB3	658531	919744	7.17%	1.044%
73	13.292	1902	1905	1907	rBV	396290	346232	2.70%	0.393%
74	13.321	1907	1910	1913	rBV	416116	335322	2.62%	0.380%
75	13.451	1929	1932	1936	rBV4	90034	163077	1.27%	0.185%
76	13.574	1950	1953	1956	rBV	172387	221642	1.73%	0.251%
77	13.716	1975	1977	1979	rVV	378781	388247	3.03%	0.441%
78	13.733	1979	1980	1983	rVV	390495	295832	2.31%	0.336%
79	13.780	1983	1988	1992	rVV3	306109	611376	4.77%	0.694%
80	13.821	1993	1995	2000	rVB4	159899	214358	1.67%	0.243%
81	13.874	2000	2004	2012	rBV2	359379	532922	4.16%	0.605%
82	13.957	2013	2018	2022	rBV2	1975388	3067685	23.92%	3.481%
83	13.992	2022	2024	2030	rVB	1698269	1544468	12.05%	1.752%
84	14.074	2035	2038	2040	rBV	327935	285500	2.23%	0.324%
85	14.127	2044	2047	2051	rBV2	184165	183909	1.43%	0.209%
86	14.186	2055	2057	2061	rVB2	112112	167353	1.31%	0.190%
87	14.351	2080	2085	2088	rVV	541618	594431	4.64%	0.674%
88	14.386	2088	2091	2095	rBV	258642	247239	1.93%	0.281%
89	14.451	2100	2102	2104	rVB2	184440	139853	1.09%	0.159%
90	14.480	2104	2107	2108	rBV	257686	232031	1.81%	0.263%
91	14.821	2161	2165	2167	rBV	205551	276372	2.16%	0.314%
92	15.010	2192	2197	2199	rBV	1285420	1778477	13.87%	2.018%
93	15.115	2211	2215	2220	rVB	315330	357504	2.79%	0.406%
94	15.310	2243	2248	2254	rBV	748630	930863	7.26%	1.056%
95	15.374	2254	2259	2265	rVV	1150240	1566163	12.21%	1.777%
96	15.433	2265	2269	2271	rVV	570950	708753	5.53%	0.804%
97	15.462	2271	2274	2280	rVB2	346501	478130	3.73%	0.543%
98	15.992	2361	2364	2369	rBV2	108095	138676	1.08%	0.157%
99	16.904	2514	2519	2526	rVB2	465712	904465	7.05%	1.026%
100	17.351	2590	2595	2607	rVB	359983	799608	6.24%	0.907%

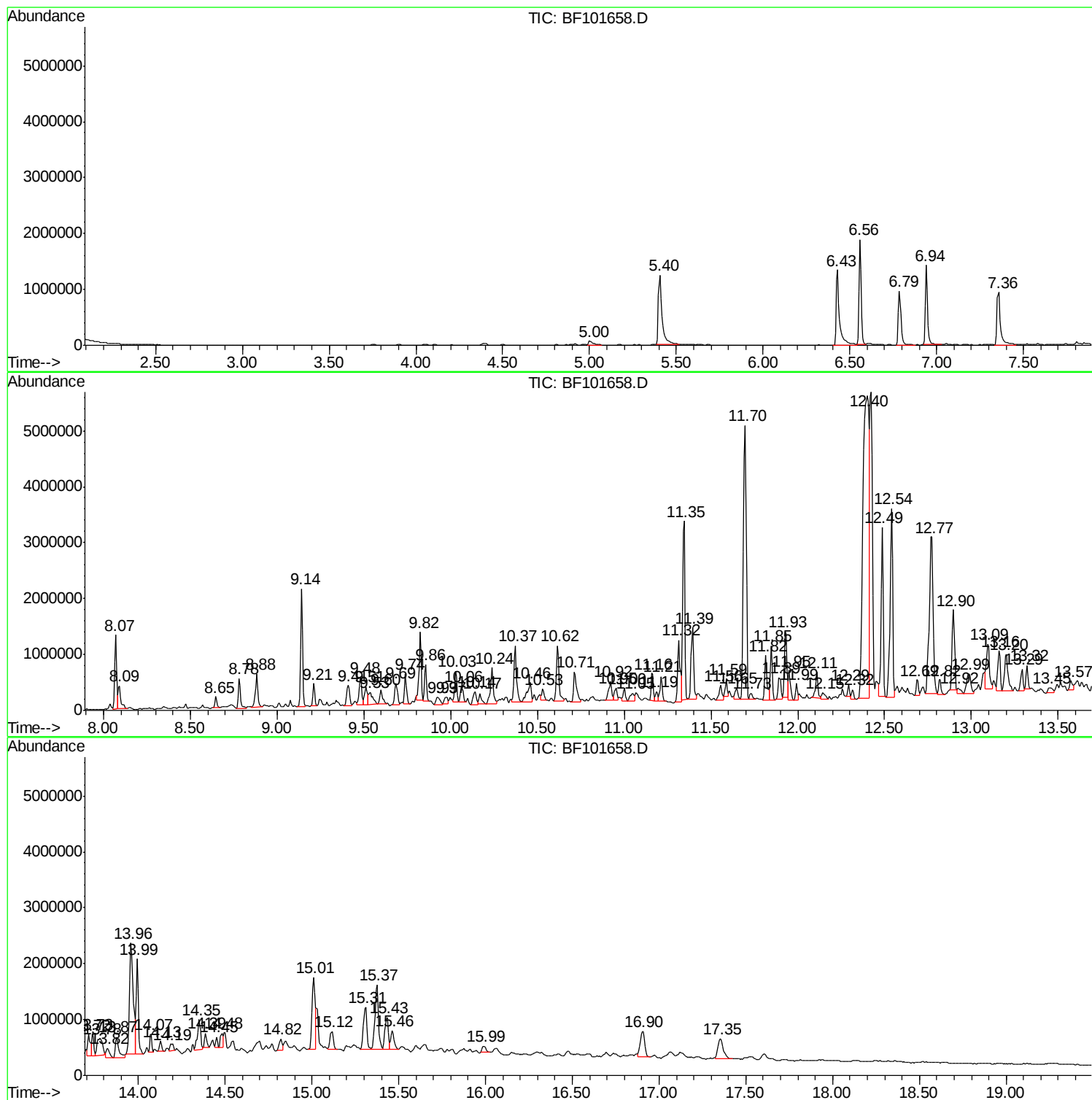
Sum of corrected areas: 88131646

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF121417.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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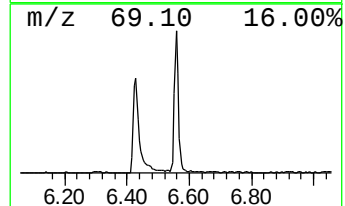
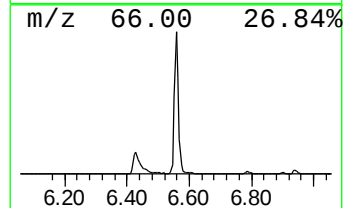
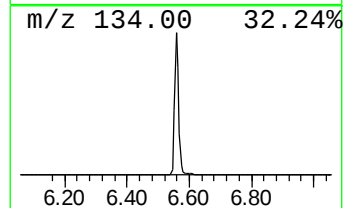
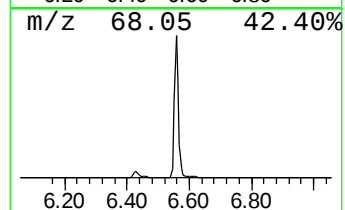
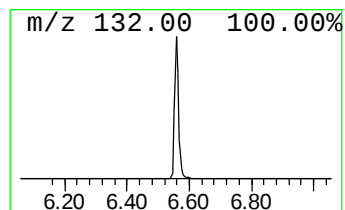
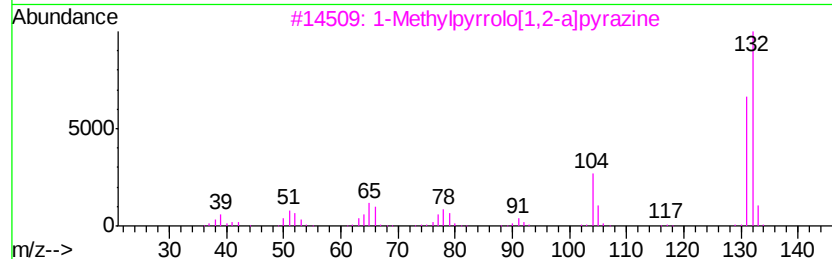
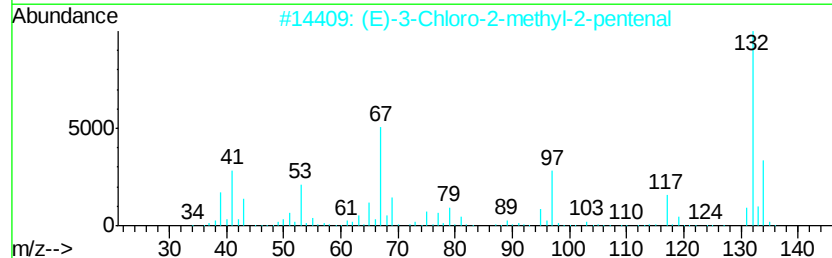
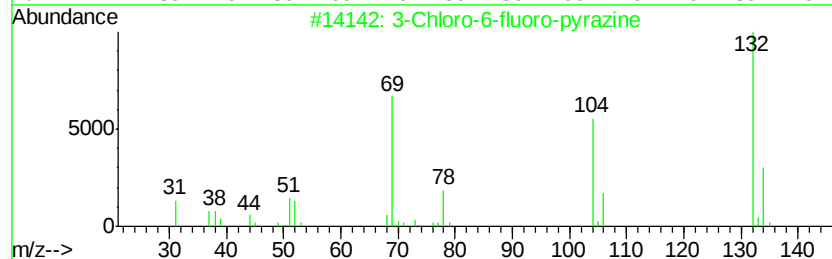
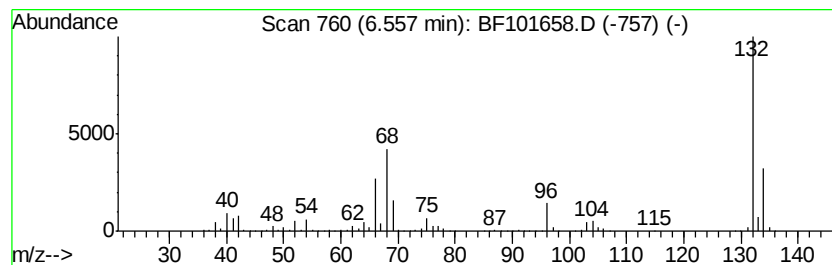
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF121417.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 unknown6.56 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.56	37.46 ng	1724290	1,4-Dichlorobenzene-d4	6.79

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
3		1-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-59-9	9
4		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9
5		1,3,5-Trifluorobenzene	132	C6H3F3	000372-38-3	9



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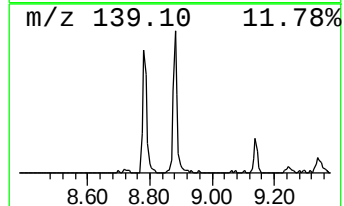
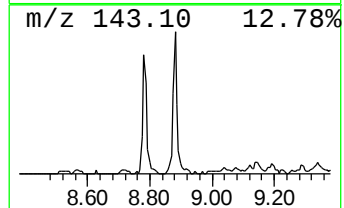
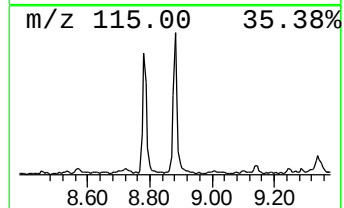
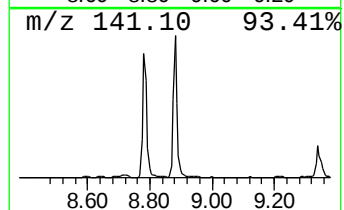
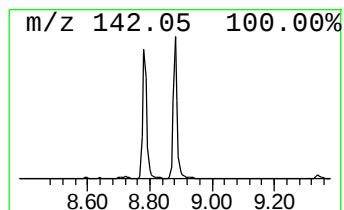
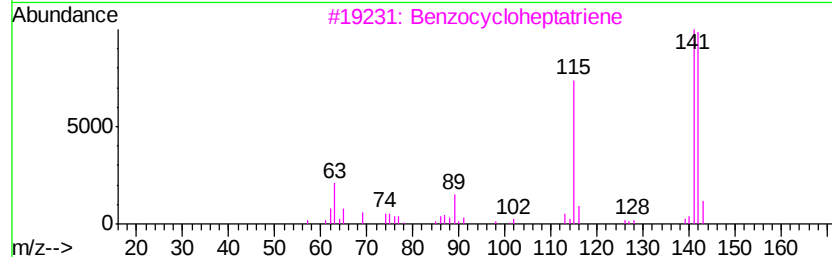
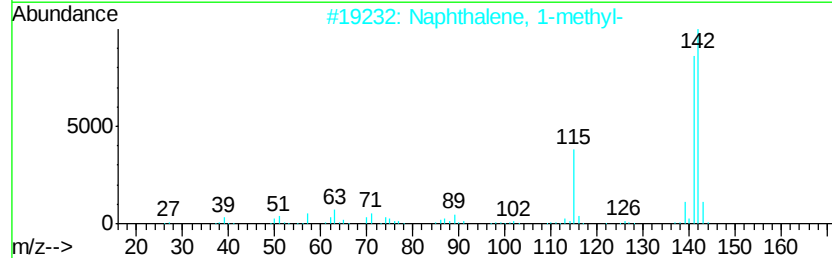
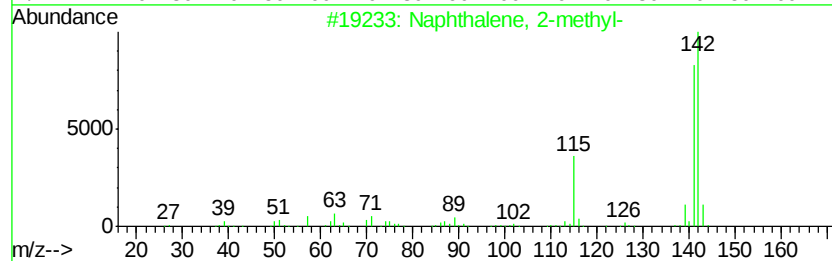
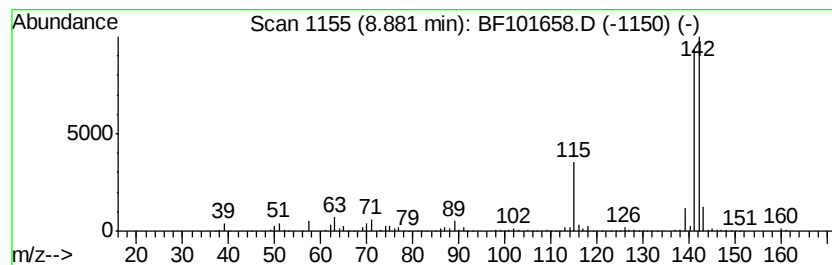
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TIC Integration Parameters: LSCINT.P

Peak Number 3 Naphthalene, 1-methyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.88	9.92 ng	565508	Naphthalene-d8	8.07

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
2		Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
3		Benzocycloheptatriene	142	C11H10	000264-09-5	94
4		1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	91
5		1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	86



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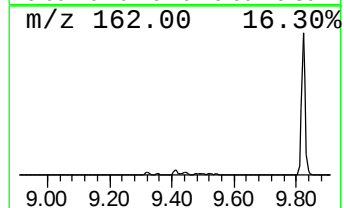
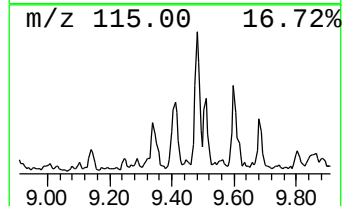
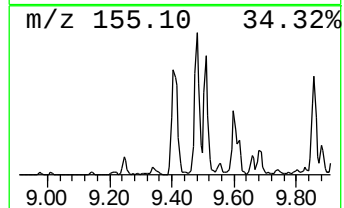
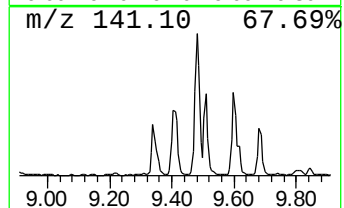
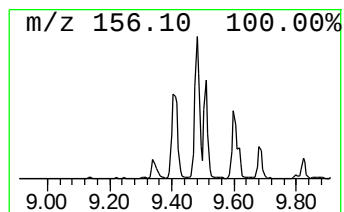
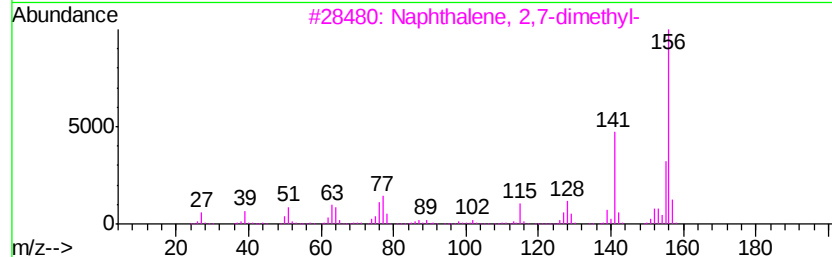
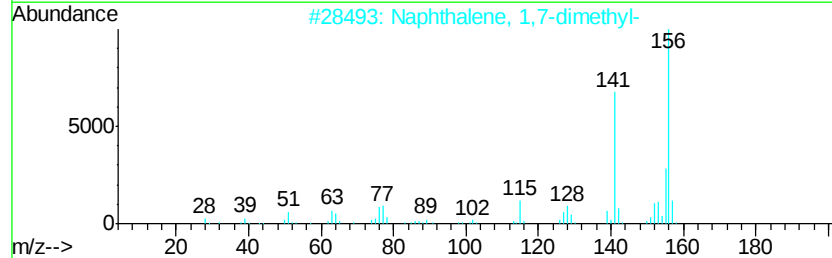
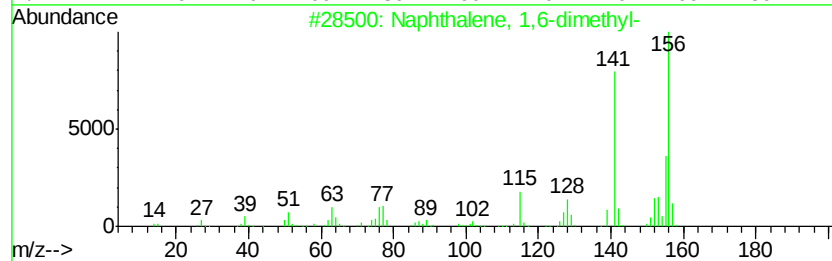
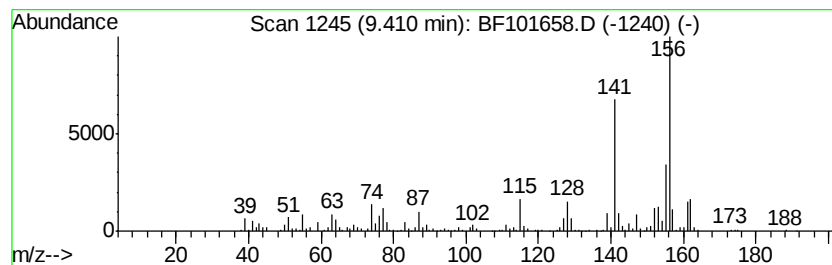
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Naphthalene, 1,6-dimethyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.41	9.46 ng	479357	Acenaphthene-d10	9.82
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 1,6-dimethyl-	156 C12H12	000575-43-9 98
2		Naphthalene, 1,7-dimethyl-	156 C12H12	000575-37-1 97
3		Naphthalene, 2,7-dimethyl-	156 C12H12	000582-16-1 97
4		Naphthalene, 2,3-dimethyl-	156 C12H12	000581-40-8 96
5		Naphthalene, 1,5-dimethyl-	156 C12H12	000571-61-9 96



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF122217\
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Operator : SJ/JU
Sample : I6683-09 2X
Misc :
ALS Vial : 10 Sample Multiplier: 1

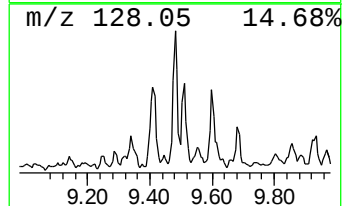
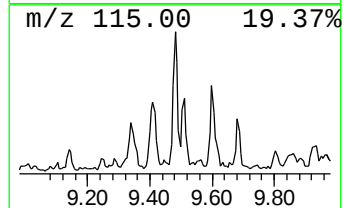
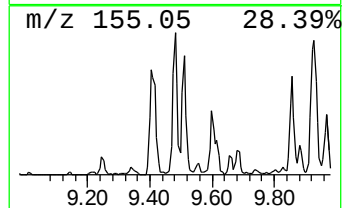
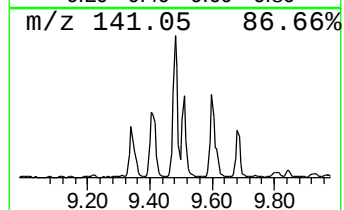
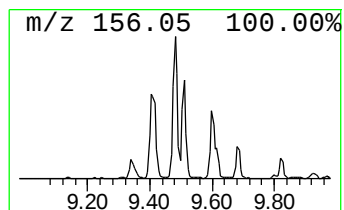
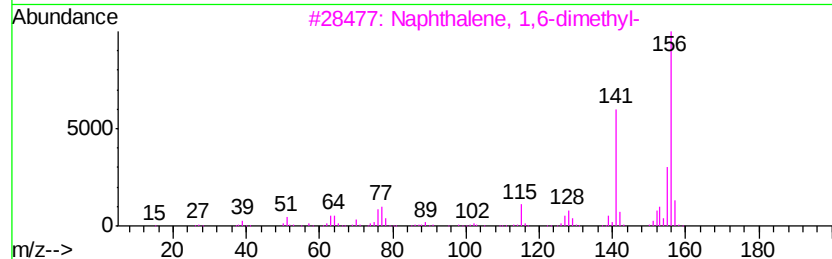
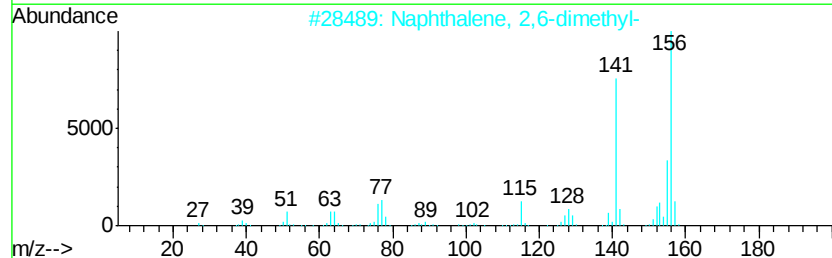
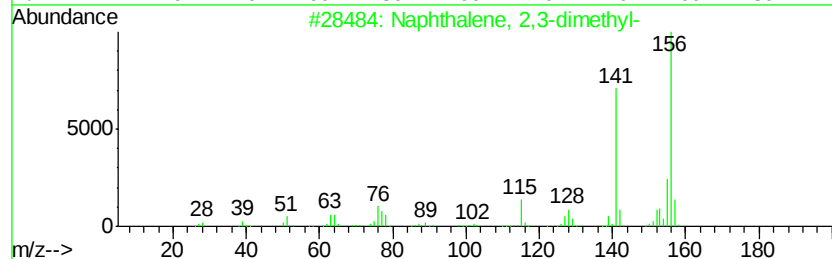
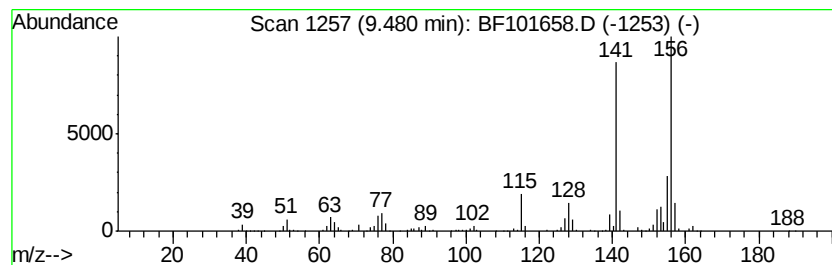
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Naphthalene, 2,3-dimethyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.48	11.17 ng	566048	Acenaphthene-d10	9.82
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 2,3-dimethyl-	156 C12H12	000581-40-8 97
2		Naphthalene, 2,6-dimethyl-	156 C12H12	000581-42-0 97
3		Naphthalene, 1,6-dimethyl-	156 C12H12	000575-43-9 97
4		Naphthalene, 1,3-dimethyl-	156 C12H12	000575-41-7 97
5		Naphthalene, 1,4-dimethyl-	156 C12H12	000571-58-4 97



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF122217\
Data File : BF101658.D
Acq On : 22 Dec 2017 17:26
Operator : SJ/JU
Sample : I6683-09 2X
Misc :
ALS Vial : 10 Sample Multiplier: 1

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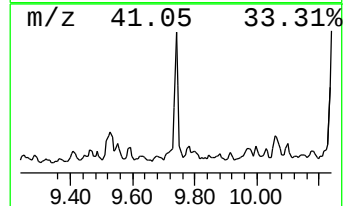
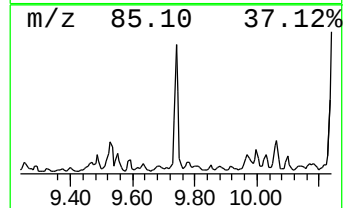
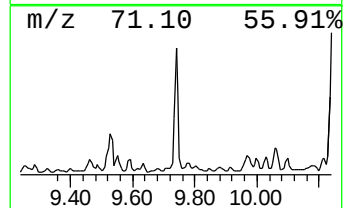
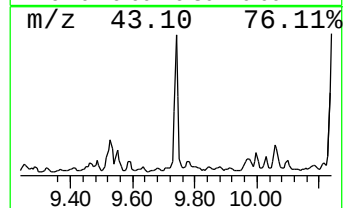
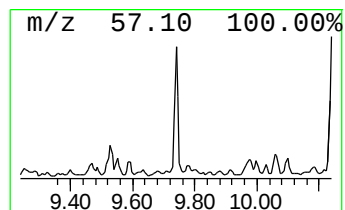
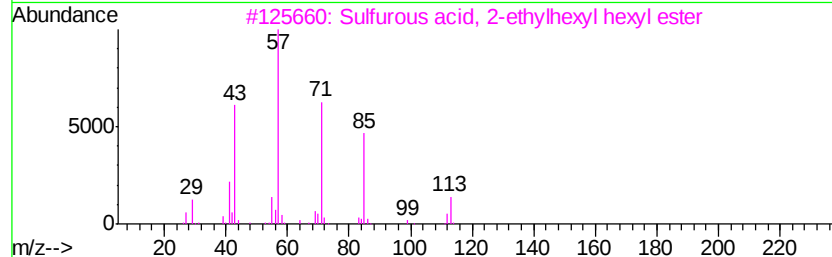
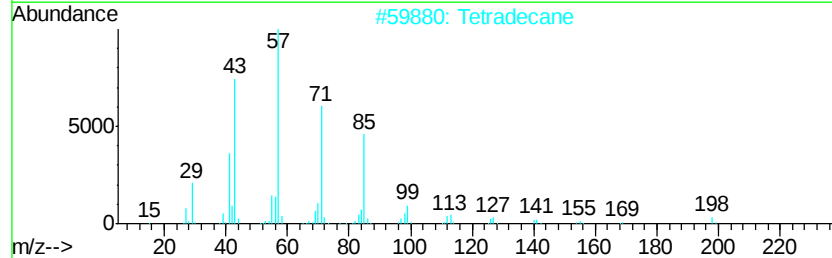
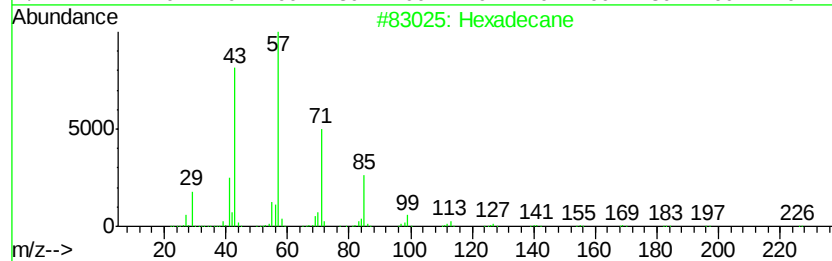
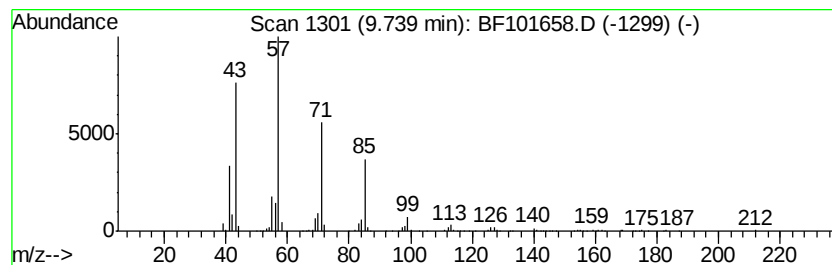
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Hexadecane Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.74	8.11 ng	410883	Acenaphthene-d10	9.82

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane		226	C16H34	000544-76-3	91
2	Tetradecane		198	C14H30	000629-59-4	90
3	Sulfurous acid, 2-ethylhexyl hex...		278	C14H30O3S	1000309-20-2	86
4	Bacchotricuneatin c		342	C20H22O5	066563-30-2	81
5	Nonane, 3,7-dimethyl-		156	C11H24	017302-32-8	72



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF122217\
Data File : BF101658.D
Acq On : 22 Dec 2017 17:26
Operator : SJ/JU
Sample : I6683-09 2X
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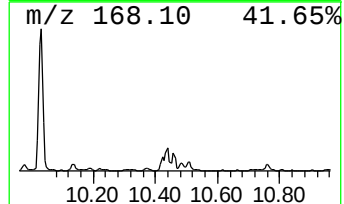
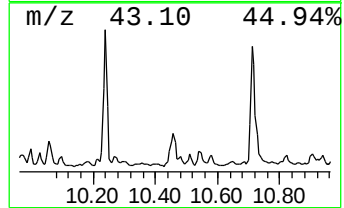
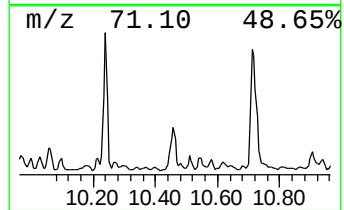
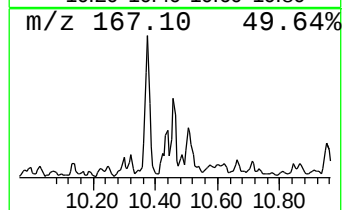
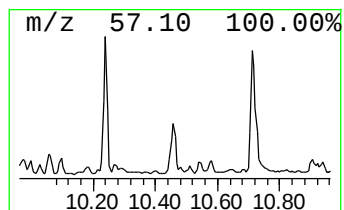
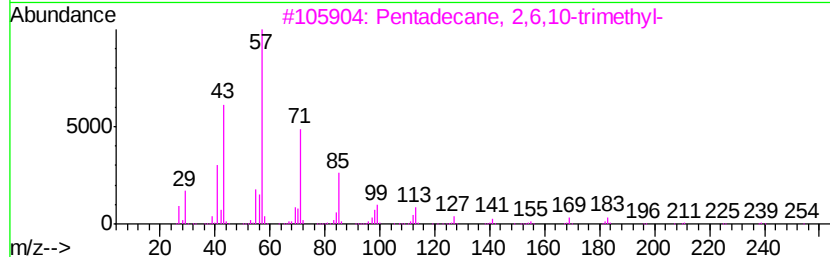
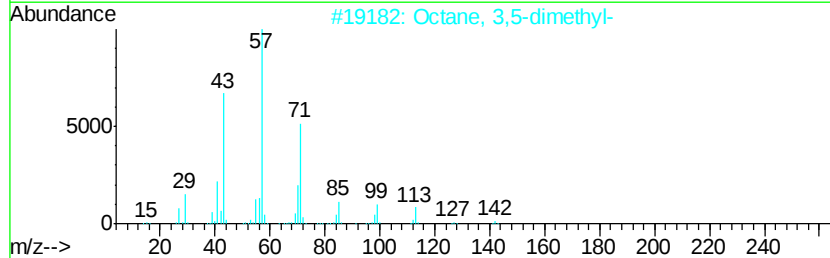
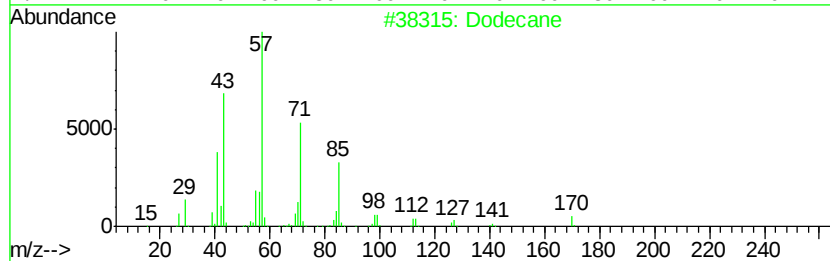
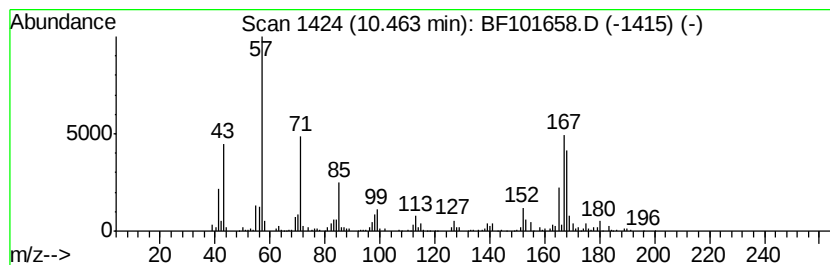
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF121417.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 unknown10.46 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.46	10.77 ng	545992	Acenaphthene-d10	9.82
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Dodecane	170 C12H26	000112-40-3	20
2	Octane, 3,5-dimethyl-	142 C10H22	015869-93-9	15
3	Pentadecane, 2,6,10-trimethyl-	254 C18H38	003892-00-0	11
4	Heptadecane, 2,6,10,15-tetramethyl-	296 C21H44	054833-48-6	11
5	Tridecane	184 C13H28	000629-50-5	11



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF122217\
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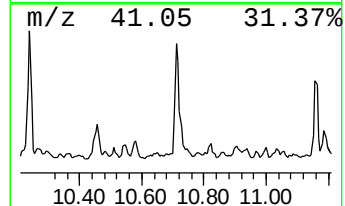
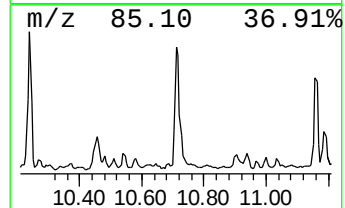
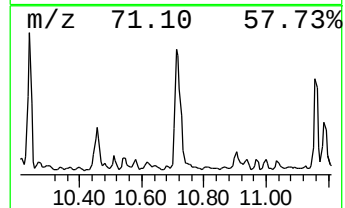
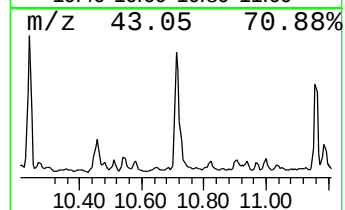
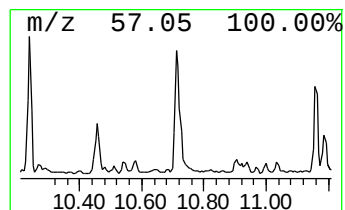
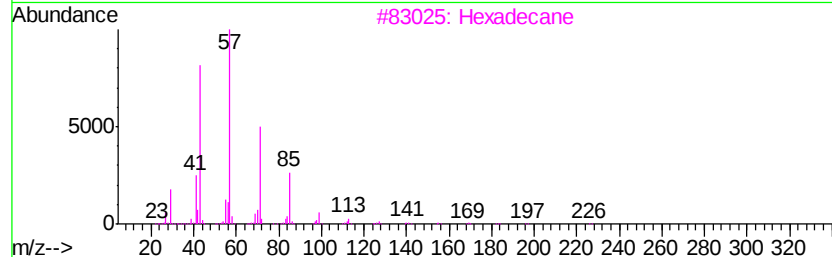
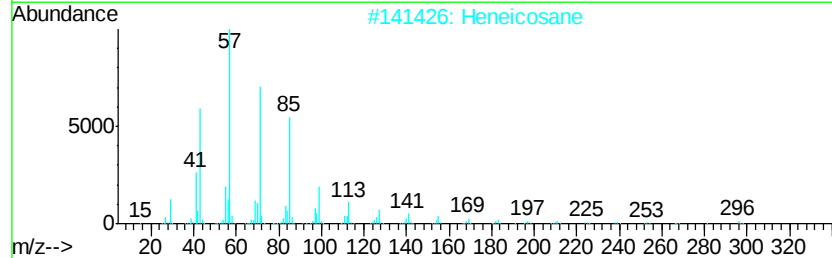
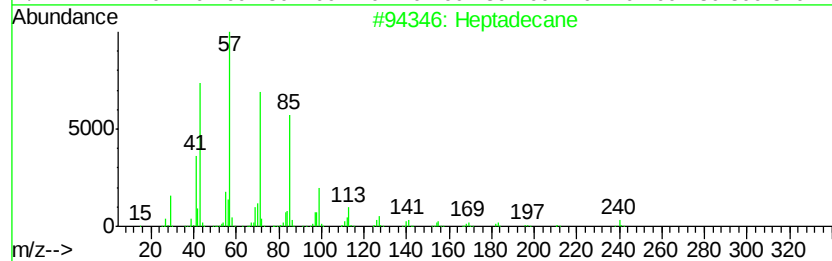
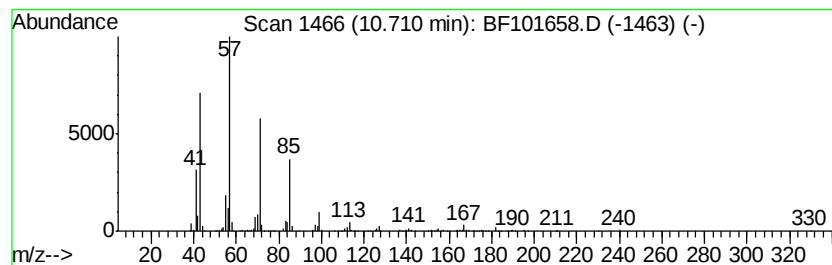
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Heptadecane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.71	12.33 ng	691187	Phenanthrene-d10	11.32

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptadecane	240	C17H36	000629-78-7	95
2		Heneicosane	296	C21H44	000629-94-7	90
3		Hexadecane	226	C16H34	000544-76-3	86
4		Dodecane, 2-methyl-6-propyl-	226	C16H34	055045-08-4	86
5		Tridecane, 1-iodo-	310	C13H27I	035599-77-0	78



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF122217\
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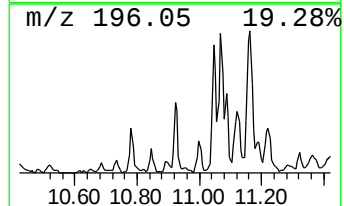
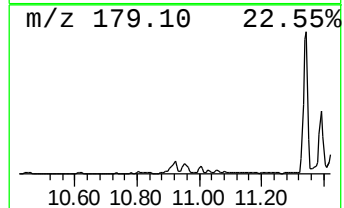
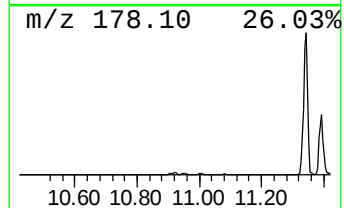
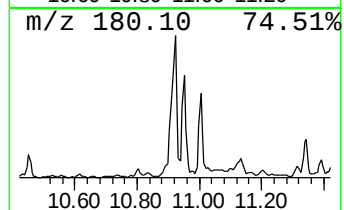
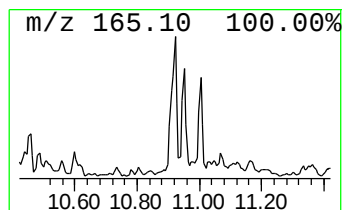
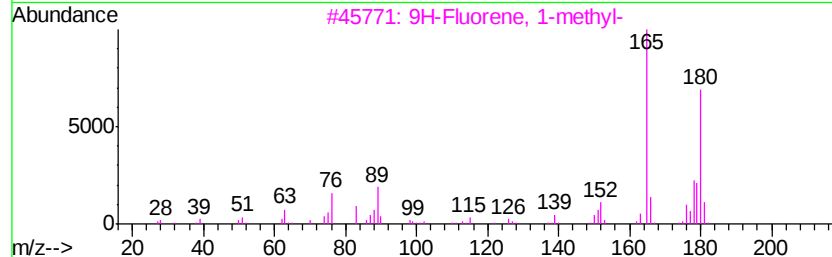
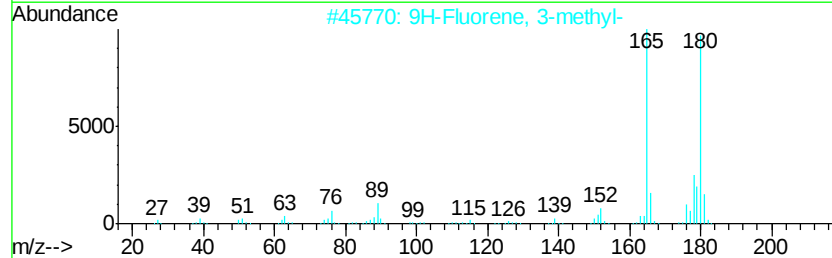
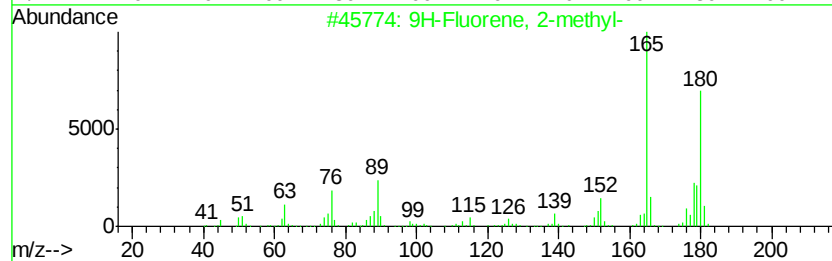
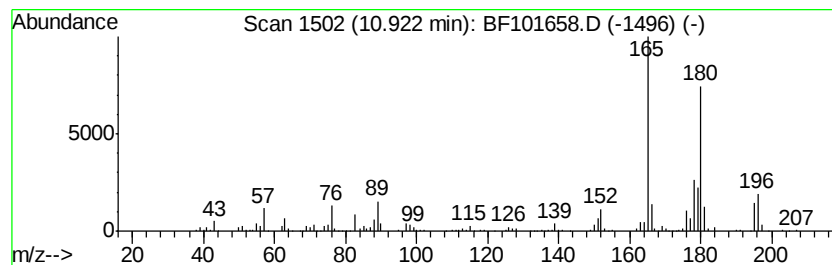
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 9H-Fluorene, 2-methyl- Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.92	8.34 ng	467437	Phenanthrene-d10	11.32

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			9H-Fluorene, 2-methyl-	180	C14H12	001430-97-3	97
2			9H-Fluorene, 3-methyl-	180	C14H12	002523-39-9	95
3			9H-Fluorene, 1-methyl-	180	C14H12	001730-37-6	93
4			9H-Fluorene, 9-methyl-	180	C14H12	002523-37-7	93
5			4a,9a-Methano-9H-fluorene	180	C14H12	019540-84-2	86



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF122217\
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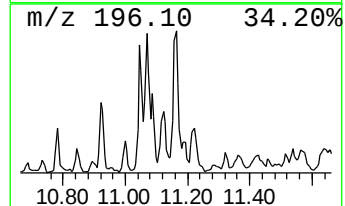
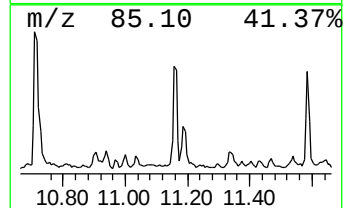
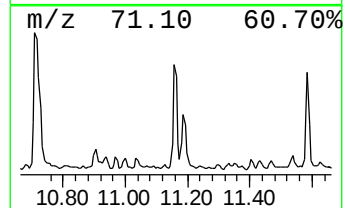
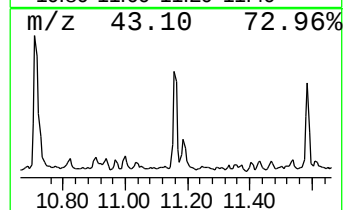
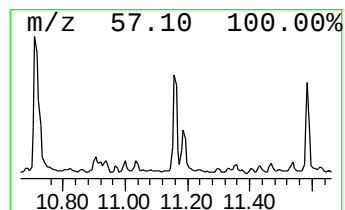
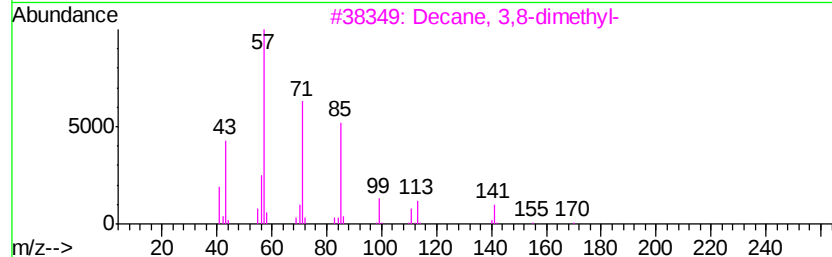
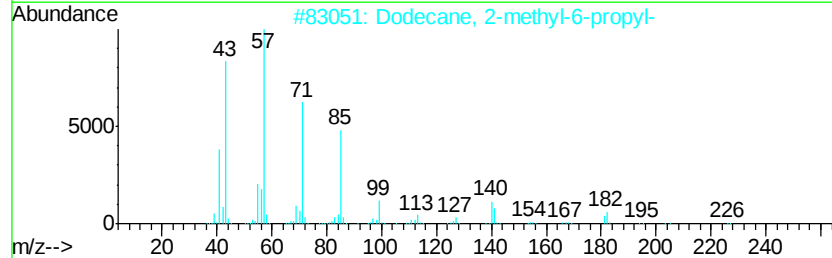
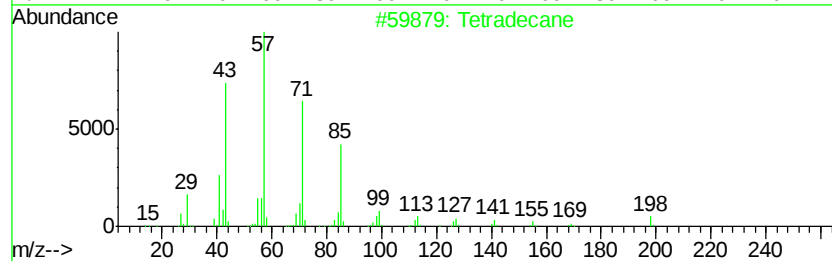
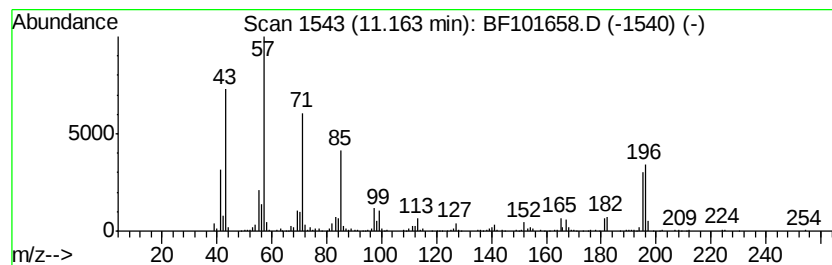
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Tetradecane Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD		R.T.
11.16	8.22 ng	460781	Phenanthrene-d10		11.32
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetradecane	198	C14H30	000629-59-4	91
2	Dodecane, 2-methyl-6-propyl-	226	C16H34	055045-08-4	58
3	Decane, 3,8-dimethyl-	170	C12H26	017312-55-9	53
4	Heneicosane	296	C21H44	000629-94-7	52
5	Undecane, 3-methyl-	170	C12H26	001002-43-3	52



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF122217\
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Sample : I6683-09 2X
Misc :
ALS Vial : 10 Sample Multiplier: 1

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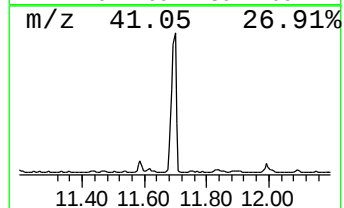
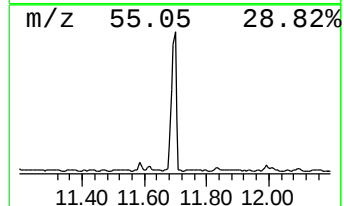
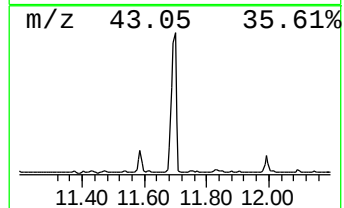
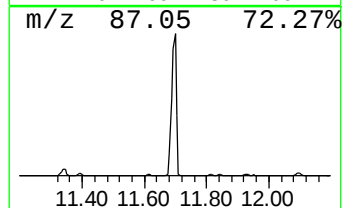
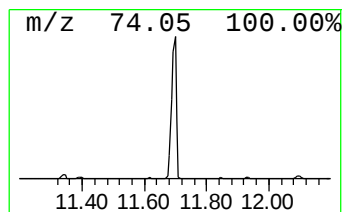
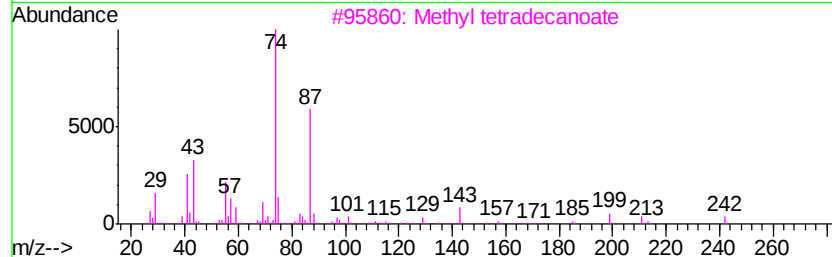
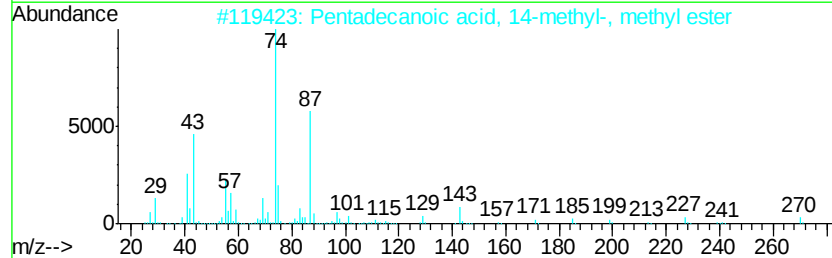
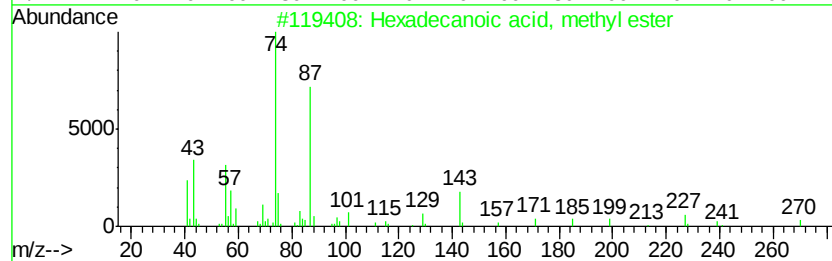
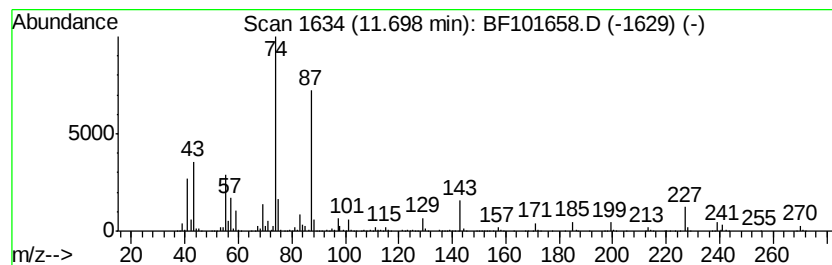
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Hexadecanoic acid, methyl e... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.70	92.87 ng	5205440	Phenanthrene-d10	11.32

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	98
2		Pentadecanoic acid, 14-methyl-, ...	270	C17H34O2	005129-60-2	96
3		Methyl tetradecanoate	242	C15H30O2	000124-10-7	91
4		Tridecanoic acid, methyl ester	228	C14H28O2	001731-88-0	89
5		Pentadecanoic acid, methyl ester	256	C16H32O2	007132-64-1	86



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF122217\
Data File : BF101658.D
Acq On : 22 Dec 2017 17:26
Operator : SJ/JU
Sample : I6683-09 2X
Misc :
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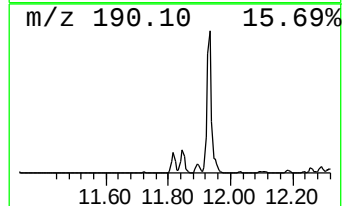
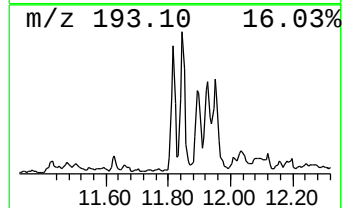
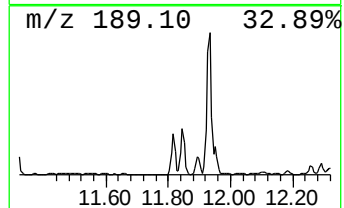
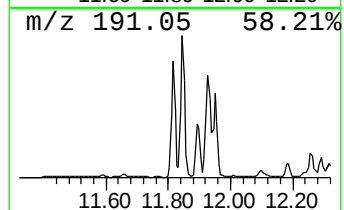
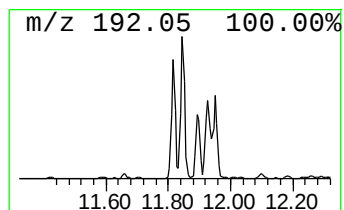
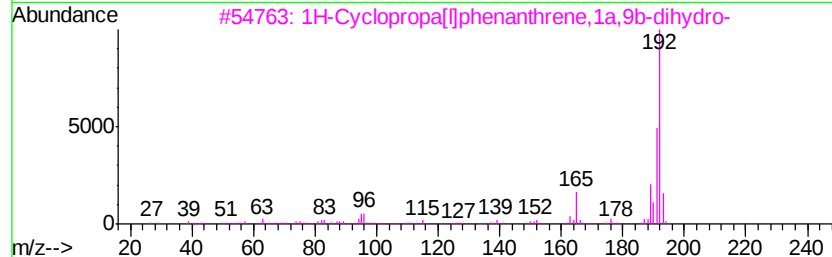
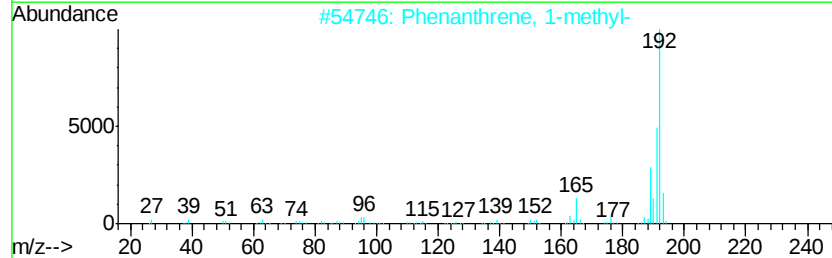
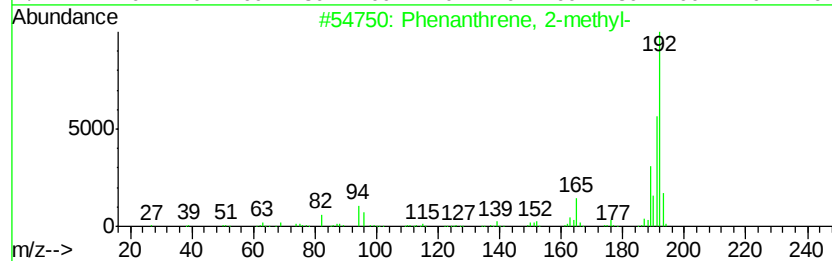
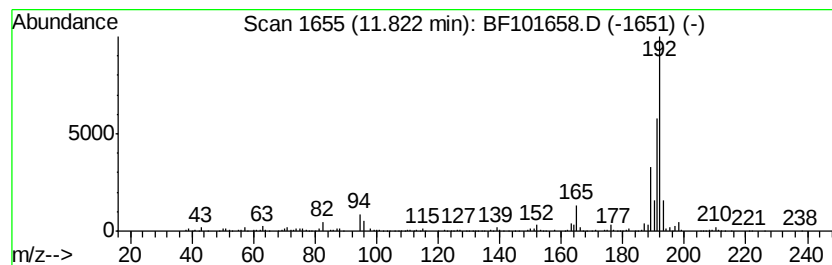
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Phenanthrene, 2-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.82	13.53 ng	758129	Phenanthrene-d10	11.32	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	98
2	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	97
3	1H-Cyclopropa[1]phenanthrene,1a,...	192	C15H12	000949-41-7	96
4	Anthracene, 2-methyl-	192	C15H12	000613-12-7	96
5	Anthracene, 9-methyl-	192	C15H12	000779-02-2	95



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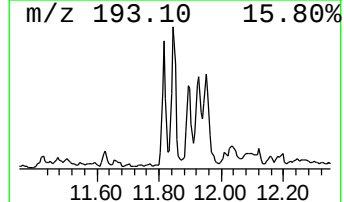
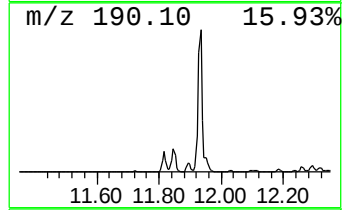
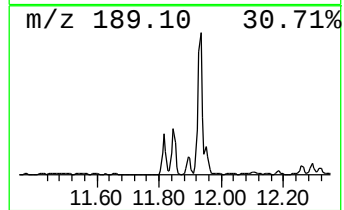
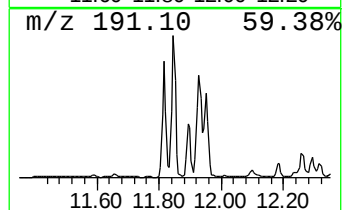
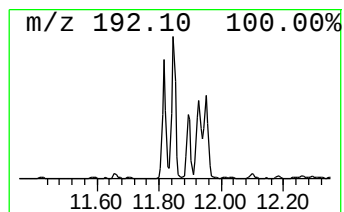
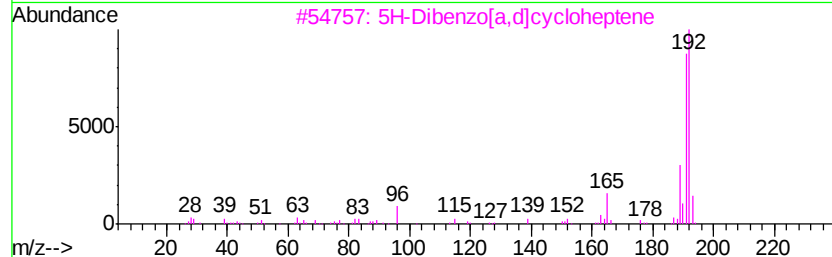
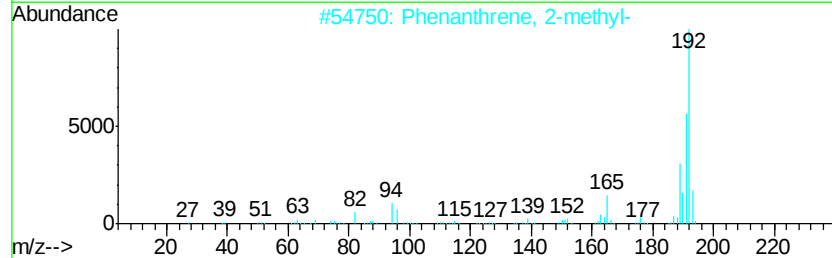
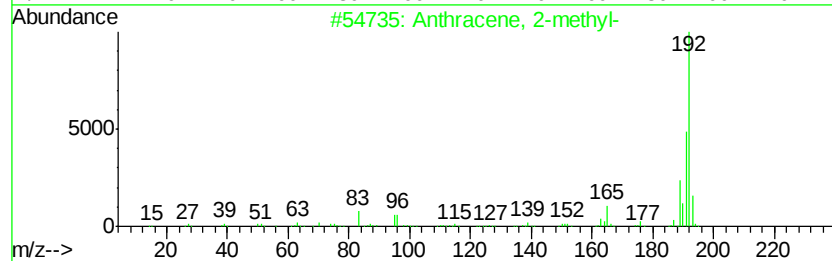
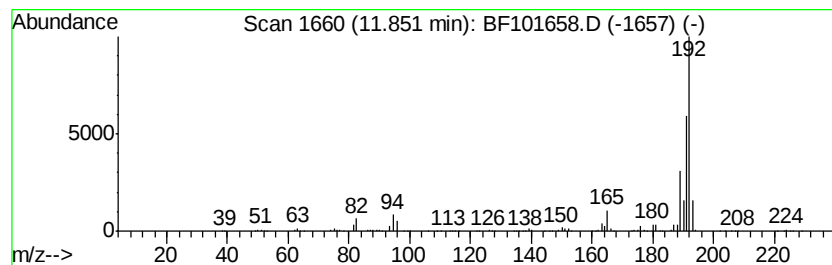
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Anthracene, 2-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.85	14.47 ng	810749	Phenanthrene-d10	11.32

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Anthracene, 2-methyl-	192	C15H12	000613-12-7	96
2		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	96
3		5H-Dibenzo[a,d]cycloheptene	192	C15H12	000256-81-5	96
4		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	96
5		1H-Cyclopropa[1]phenanthrene, 1a, ...	192	C15H12	000949-41-7	95



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Misc :
ALS Vial : 10 Sample Multiplier: 1

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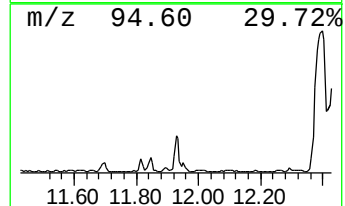
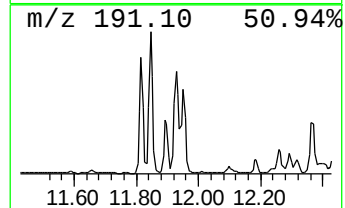
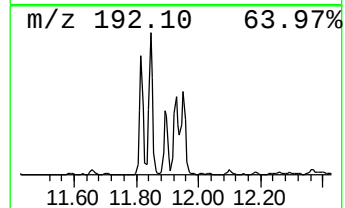
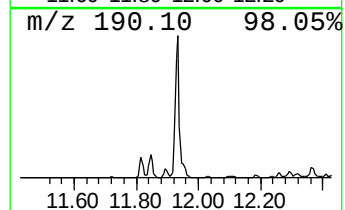
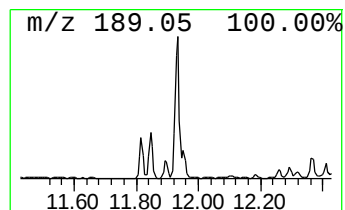
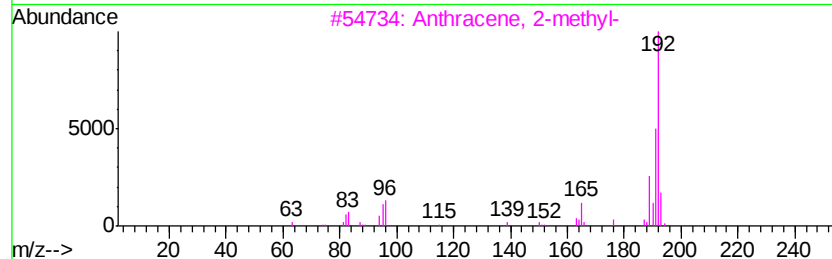
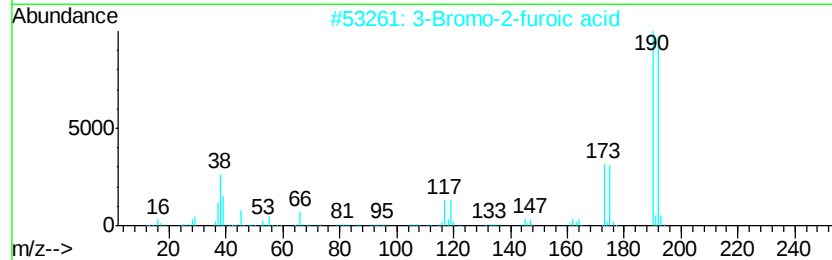
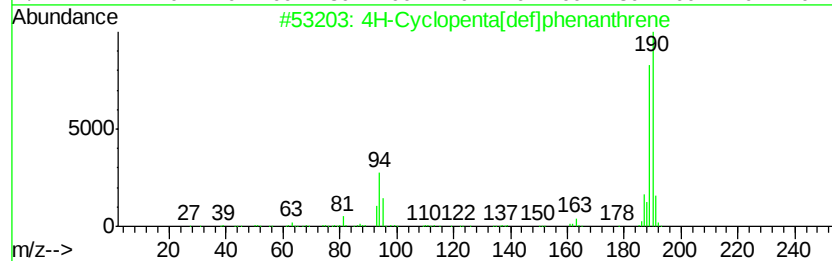
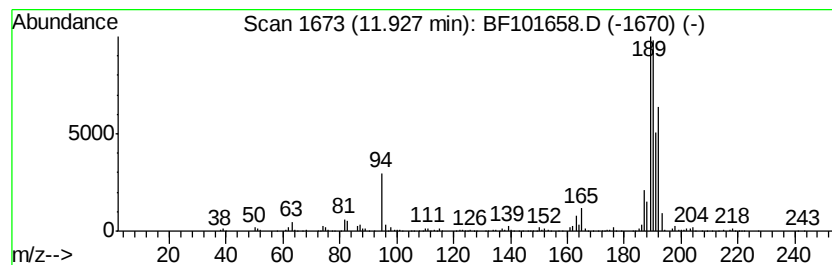
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 4H-Cyclopenta[def]phenanthrene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.93	23.64 ng	1325220	Phenanthrene-d10	11.32

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	70
2		3-Bromo-2-furoic acid	190	C5H3BrO3	014903-90-3	27
3		Anthracene, 2-methyl-	192	C15H12	000613-12-7	18
4		8,9-Dihydrocyclopenta[def]phenan...	192	C15H12	027410-55-5	18
5		1a,9b-Dihydro-1H-cyclopropa[a]an...	192	C15H12	006109-24-6	18



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF122217\
Data File : BF101658.D
Acq On : 22 Dec 2017 17:26
Operator : SJ/JU
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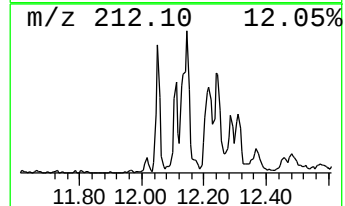
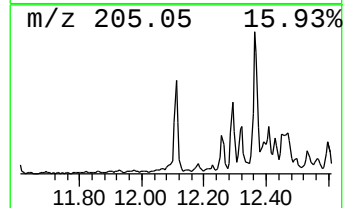
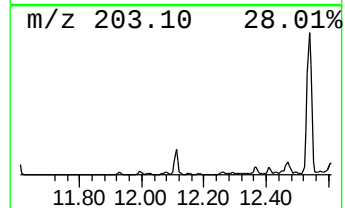
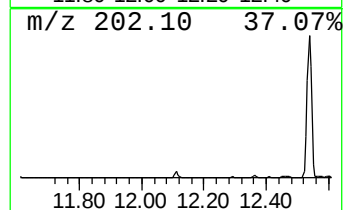
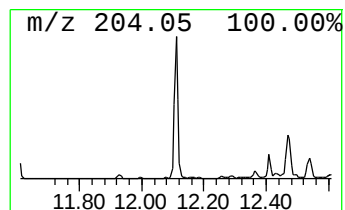
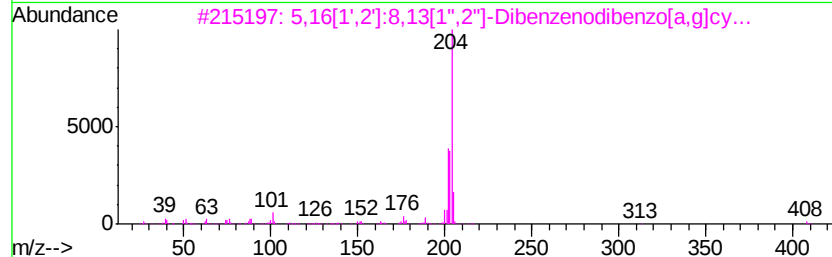
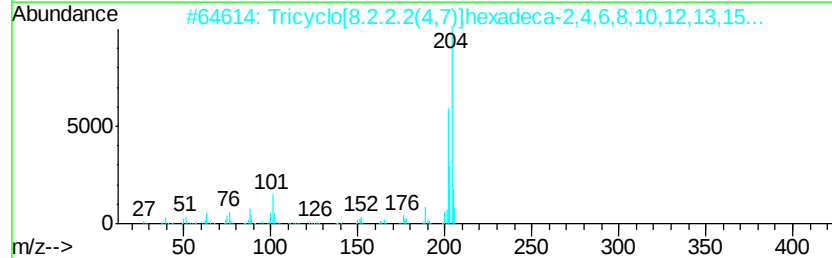
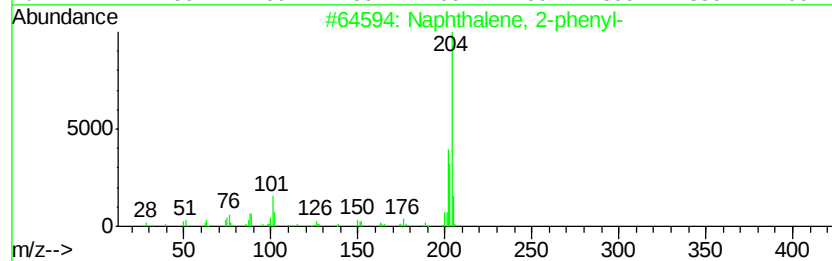
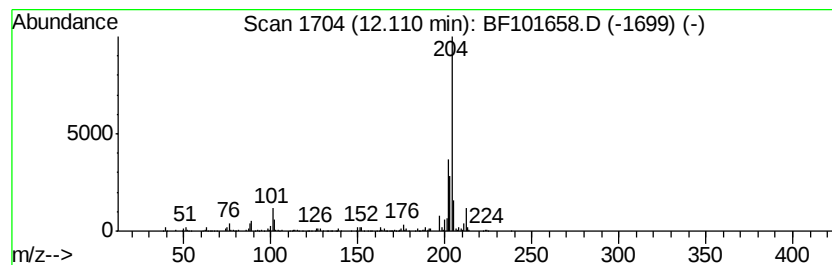
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 Naphthalene, 2-phenyl- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.11	8.47 ng	474656	Phenanthrene-d10	11.32

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	96
2		Tricyclo[8.2.2.2(4,7)]hexadeca-2...	204	C16H12	006572-60-7	64
3		5,16[1',2']:8,13[1'',2'']-Dibenz...	408	C32H24	005672-97-9	59
4		9,10-Bis(bromomethyl)anthracene	362	C16H12Br2	034373-96-1	58
5		Acephenanthrylene, 4,5-dihydro-	204	C16H12	006232-48-0	53



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF122217\
Data File : BF101658.D
Acq On : 22 Dec 2017 17:26
Operator : SJ/JU
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Misc :
ALS Vial : 10 Sample Multiplier: 1

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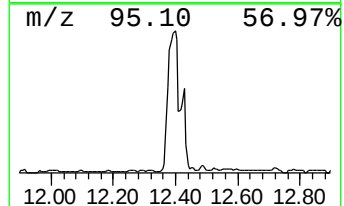
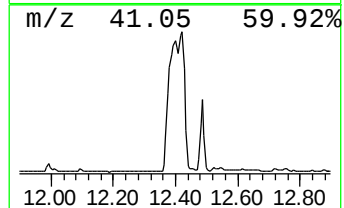
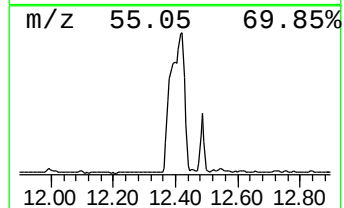
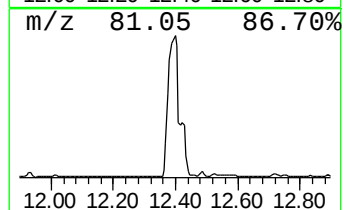
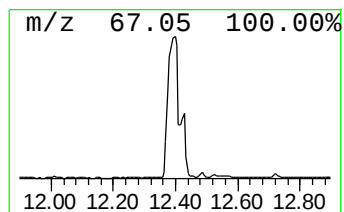
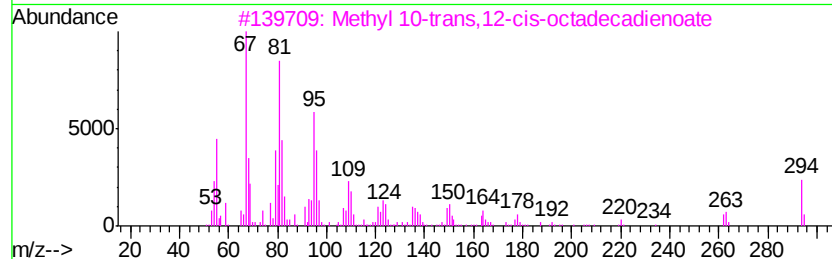
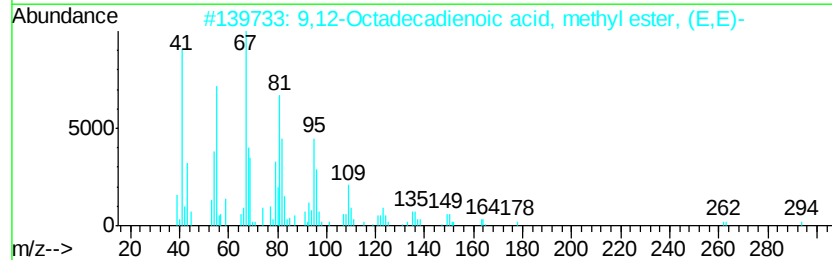
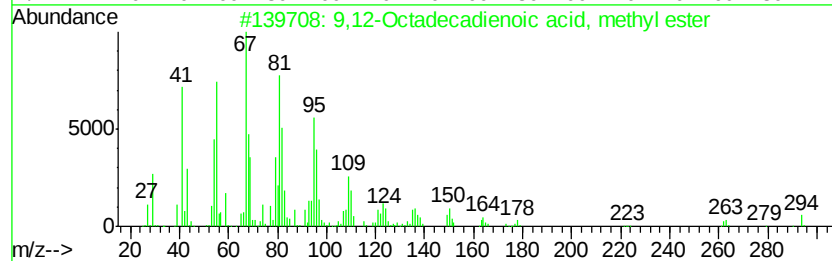
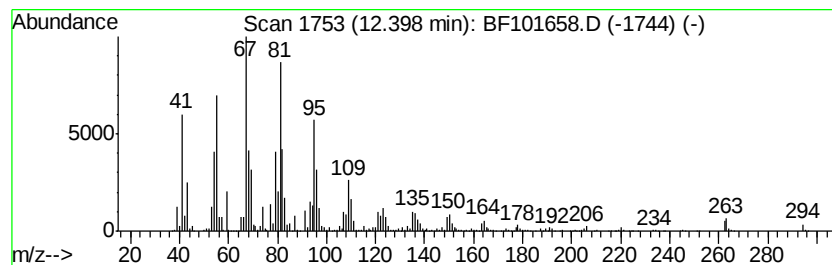
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 9,12-Octadecadienoic acid, ... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.40	228.77 ng	12822100	Phenanthrene-d10	11.32

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	9,12-Octadecadienoic acid, methy...	294	C19H34O2	002462-85-3	99
2		9,12-Octadecadienoic acid, methy...	294	C19H34O2	002566-97-4	99
3		Methyl 10-trans,12-cis-octadecad...	294	C19H34O2	1000336-44-2	99
4		9,15-Octadecadienoic acid, methy...	294	C19H34O2	017309-05-6	99
5		9,12-Octadecadienoic acid (Z,Z)-...	294	C19H34O2	000112-63-0	98



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF122217\
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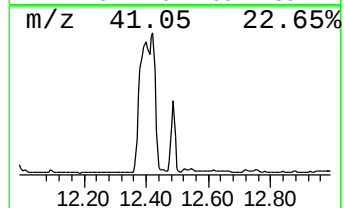
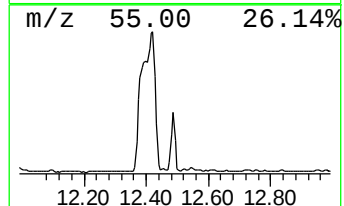
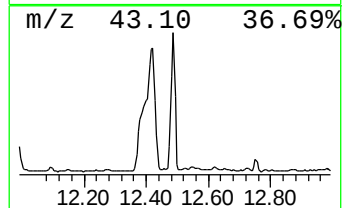
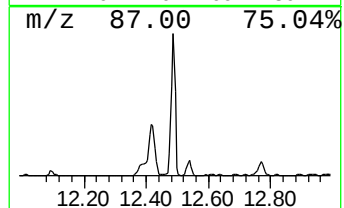
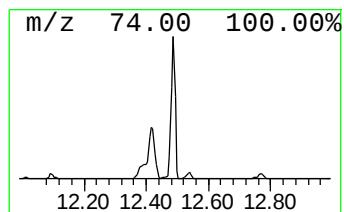
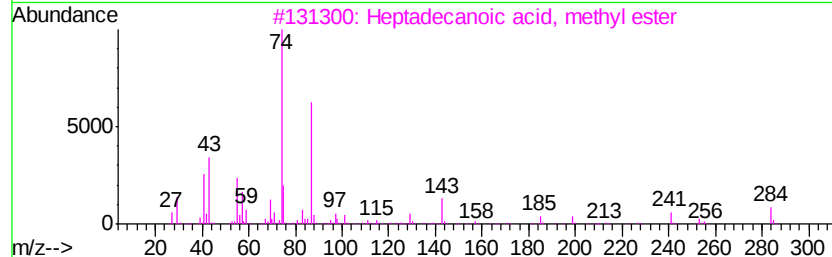
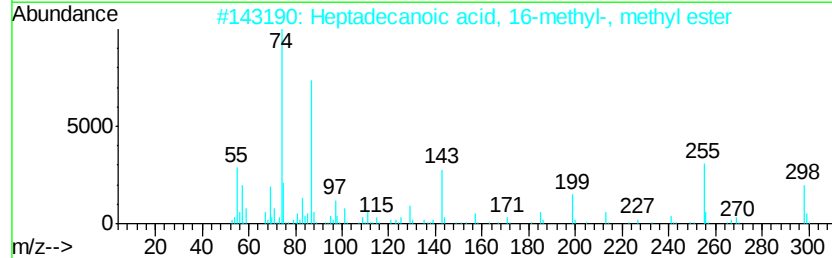
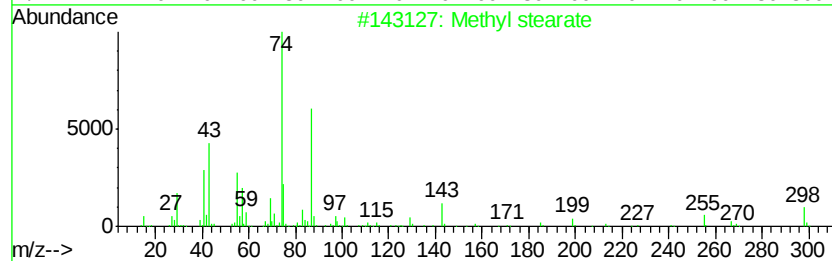
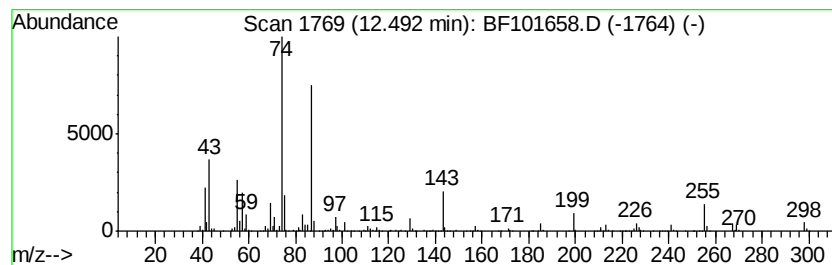
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 Methyl stearate Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.49	49.29 ng	2762390	Phenanthrene-d10	11.32
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Methyl stearate	298 C19H38O2	000112-61-8 97
2		Heptadecanoic acid, 16-methyl-, ...	298 C19H38O2	005129-61-3 91
3		Heptadecanoic acid, methyl ester	284 C18H36O2	001731-92-6 91
4		Hexadecanoic acid, methyl ester	270 C17H34O2	000112-39-0 90
5		Methyl tetradecanoate	242 C15H30O2	000124-10-7 90



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Misc :
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ClientSampleId :
EO-02-122017-B

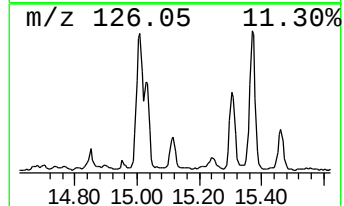
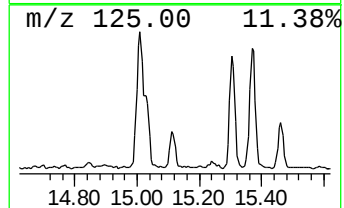
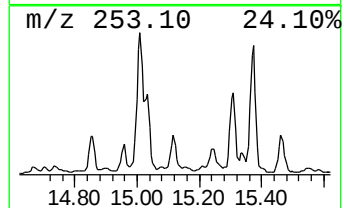
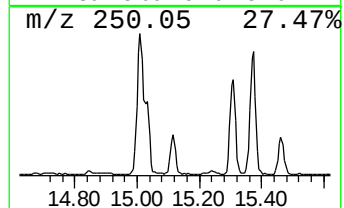
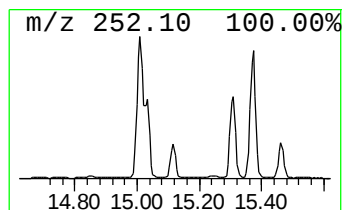
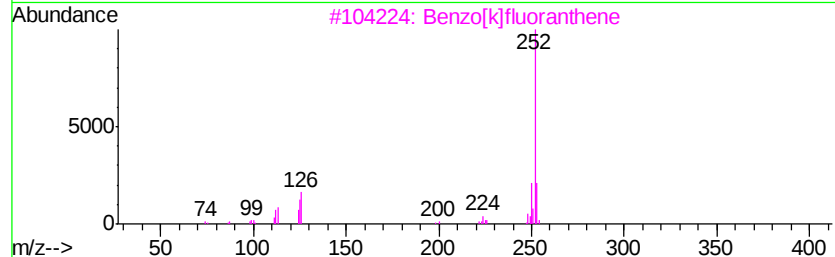
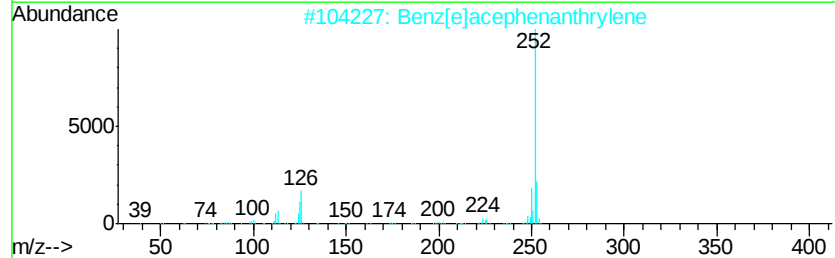
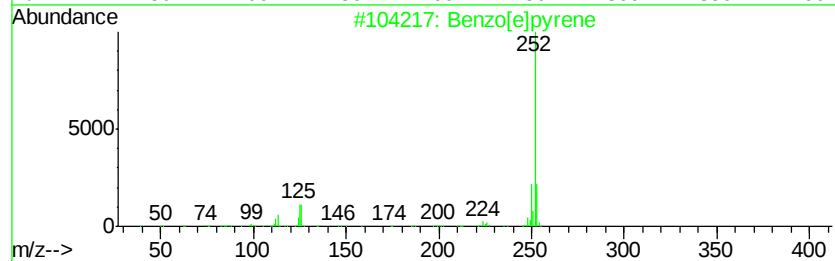
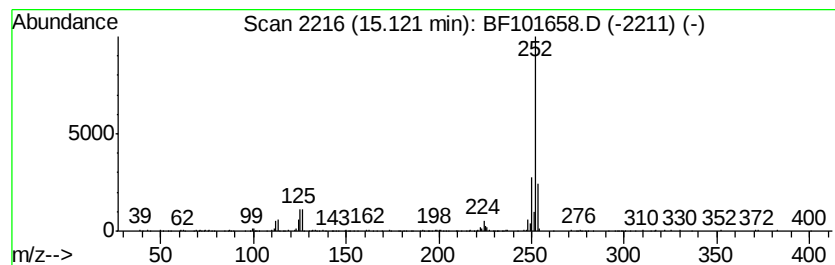
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF121417.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 19 Benzo[e]pyrene Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.12	10.09 ng	357504	Perylene-d12	15.43

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzo[e]pyrene	252	C20H12	000192-97-2	98
2		Benz[e]acephenanthrylene	252	C20H12	000205-99-2	98
3		Benzo[k]fluoranthene	252	C20H12	000207-08-9	93
4		Benzo[a]pyrene	252	C20H12	000050-32-8	93
5		Perylene	252	C20H12	000198-55-0	87



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF122217\
 Data File : BF101658.D
 Acq On : 22 Dec 2017 17:26
 Operator : SJ/JU
 Sample : I6683-09 2X
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EO-02-122017-B

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF121417.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
unknown6.56	6.56	37.5	ng	1724290	1	6.79	920636	20.0
Naphthalene, 1-me...	8.88	9.9	ng	565508	2	8.07	1139610	20.0
Naphthalene, 1,6-...	9.41	9.5	ng	479357	3	9.82	1013860	20.0
Naphthalene, 2,3-...	9.48	11.2	ng	566048	3	9.82	1013860	20.0
Hexadecane	9.74	8.1	ng	410883	3	9.82	1013860	20.0
unknown10.46	10.46	10.8	ng	545992	3	9.82	1013860	20.0
Heptadecane	10.71	12.3	ng	691187	4	11.32	1120970	20.0
9H-Fluorene, 2-me...	10.92	8.3	ng	467437	4	11.32	1120970	20.0
Tetradecane	11.16	8.2	ng	460781	4	11.32	1120970	20.0
Hexadecanoic acid...	11.70	92.9	ng	5205440	4	11.32	1120970	20.0
Phenanthrene, 2-m...	11.82	13.5	ng	758129	4	11.32	1120970	20.0
Anthracene, 2-met...	11.85	14.5	ng	810749	4	11.32	1120970	20.0
4H-Cyclopenta[def...	11.93	23.6	ng	1325220	4	11.32	1120970	20.0
Naphthalene, 2-ph...	12.11	8.5	ng	474656	4	11.32	1120970	20.0
9,12-Octadecadien...	12.40	228.8	ng	12822100	4	11.32	1120970	20.0
Methyl stearate	12.49	49.3	ng	2762390	4	11.32	1120970	20.0
Benzo[e]pyrene	15.12	10.1	ng	357504	6	15.43	708753	20.0