

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF070317\
 Data File : BF096441.D
 Acq On : 3 Jul 2017 15:14
 Operator : SJ/MA
 Sample : I3930-08 2X
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 B3-0-2

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:\HPCHEM1\BNA_F\METHODS\8270-BF070117.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	3.681	264	271	279	rBV3	27889	54579	4.60%	0.252%
2	4.328	374	381	388	rBV4	29780	55131	4.65%	0.254%
3	4.892	474	477	485	rVV	147320	188595	15.89%	0.870%
4	5.292	542	545	547	rBV	68835	68077	5.74%	0.314%
5	5.328	547	551	555	rVB	1316121	1186569	100.00%	5.474%
6	5.557	584	590	599	rVB3	75940	113923	9.60%	0.526%
7	5.639	599	604	607	rBV	47245	55860	4.71%	0.258%
8	5.881	642	645	650	rVB3	47094	52446	4.42%	0.242%
9	5.981	659	662	668	rVB2	56223	60672	5.11%	0.280%
10	6.181	689	696	699	rBV3	29810	51242	4.32%	0.236%
11	6.263	703	710	716	rVB3	45027	92599	7.80%	0.427%
12	6.328	716	721	722	rBV3	40539	42073	3.55%	0.194%
13	6.357	722	726	731	rVV	1145943	1058121	89.17%	4.881%
14	6.492	745	749	752	rBV	1146516	1062027	89.50%	4.899%
15	6.551	756	759	761	rBV	77178	60388	5.09%	0.279%
16	6.575	761	763	765	rVB	73603	50643	4.27%	0.234%
17	6.722	784	788	791	rVV	929328	746604	62.92%	3.444%
18	6.751	791	793	796	rVB2	51949	46711	3.94%	0.215%
19	6.786	796	799	804	rBV3	43630	45778	3.86%	0.211%
20	6.881	811	815	819	rVB	924336	823045	69.36%	3.797%
21	7.051	840	844	846	rBV2	49387	48589	4.09%	0.224%
22	7.122	853	856	860	rBV	56315	59404	5.01%	0.274%
23	7.281	876	883	886	rBV	784128	738405	62.23%	3.406%
24	7.333	889	892	895	rVB	134953	112328	9.47%	0.518%
25	7.375	895	899	903	rBV4	48296	57949	4.88%	0.267%
26	7.533	922	926	932	rVB4	29428	50331	4.24%	0.232%
27	8.004	1002	1006	1008	rBV	1279276	1056636	89.05%	4.874%
28	8.028	1008	1010	1012	rVB	209430	150026	12.64%	0.692%
29	8.086	1017	1020	1022	rVB	53193	44209	3.73%	0.204%
30	8.398	1070	1073	1075	rVV2	39081	41899	3.53%	0.193%
31	8.445	1078	1081	1084	rVV	120472	107784	9.08%	0.497%
32	8.610	1107	1109	1112	rVV	165363	124626	10.50%	0.575%
33	8.716	1121	1127	1130	rVB	279064	251824	21.22%	1.162%
34	8.816	1141	1144	1147	rVV	198267	156802	13.21%	0.723%

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Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF070117.M
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35	9.039	1180	1182	1185	rVB	76879	69421	5.85%	0.320%
36	9.080	1185	1189	1192	rBV	1486002	1152228	97.11%	5.315%
37	9.175	1202	1205	1210	rBV2	238287	208359	17.56%	0.961%
38	9.275	1219	1222	1227	rVB	42759	56568	4.77%	0.261%
39	9.339	1227	1233	1238	rBV2	101480	150628	12.69%	0.695%
40	9.416	1243	1246	1248	rBV	141800	118319	9.97%	0.546%
41	9.445	1248	1251	1253	rVB	105785	86773	7.31%	0.400%
42	9.475	1253	1256	1258	rBV	175914	146251	12.33%	0.675%
43	9.498	1258	1260	1262	rVB	99767	69308	5.84%	0.320%
44	9.533	1264	1266	1268	rBV	64920	46646	3.93%	0.215%
45	9.616	1276	1280	1284	rBV	208447	195918	16.51%	0.904%
46	9.663	1286	1288	1292	rVB4	42644	63431	5.35%	0.293%
47	9.704	1292	1295	1298	rVB	203326	158601	13.37%	0.732%
48	9.757	1298	1304	1307	rBV	1127930	988033	83.27%	4.558%
49	9.863	1318	1322	1326	rBV2	35591	58213	4.91%	0.269%
50	9.957	1335	1338	1342	rVB	161513	163309	13.76%	0.753%
51	9.998	1342	1345	1347	rBV	56565	49599	4.18%	0.229%
52	10.027	1347	1350	1354	rVB4	55286	60700	5.12%	0.280%
53	10.074	1354	1358	1360	rBV2	56592	67388	5.68%	0.311%
54	10.169	1370	1374	1377	rBV3	67107	88280	7.44%	0.407%
55	10.204	1377	1380	1383	rVB	192728	153059	12.90%	0.706%
56	10.298	1393	1396	1398	rVV2	78219	60670	5.11%	0.280%
57	10.422	1413	1417	1420	rBV3	87390	96341	8.12%	0.444%
58	10.457	1420	1423	1425	rVV2	64928	73730	6.21%	0.340%
59	10.539	1433	1437	1441	rVV	642409	578176	48.73%	2.667%
60	10.586	1441	1445	1449	rVB5	35962	47071	3.97%	0.217%
61	10.674	1456	1460	1461	rBV	251057	228531	19.26%	1.054%
62	10.686	1461	1462	1466	rVB	217755	178327	15.03%	0.823%
63	10.980	1510	1512	1513	rVV	46549	41495	3.50%	0.191%
64	11.016	1517	1518	1520	rVV	61491	42011	3.54%	0.194%
65	11.039	1520	1522	1526	rVV	96153	86086	7.26%	0.397%
66	11.086	1527	1530	1533	rVV3	44014	59837	5.04%	0.276%
67	11.121	1533	1536	1539	rVV2	208263	187269	15.78%	0.864%
68	11.151	1539	1541	1546	rVB	73944	72139	6.08%	0.333%
69	11.233	1552	1555	1557	rBV	812717	685696	57.79%	3.163%
70	11.257	1557	1559	1562	rVB	325472	250373	21.10%	1.155%
71	11.310	1565	1568	1571	rVB	235547	179979	15.17%	0.830%

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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	11.463	1591	1594	1597	rBV	59811	60441	5.09%	0.279%
73	11.545	1605	1608	1611	rBV	137815	119122	10.04%	0.550%
74	11.739	1638	1641	1643	rBV	61301	59572	5.02%	0.275%
75	11.768	1643	1646	1650	rVB2	181081	216023	18.21%	0.997%
76	11.810	1650	1653	1656	rVB2	71446	62636	5.28%	0.289%
77	11.845	1656	1659	1661	rBV	116991	117874	9.93%	0.544%
78	11.868	1661	1663	1669	rVB	74414	65744	5.54%	0.303%
79	11.951	1675	1677	1683	rVB	133514	140068	11.80%	0.646%
80	12.033	1688	1691	1692	rVV	67789	60382	5.09%	0.279%
81	12.051	1692	1694	1699	rVB2	79314	70580	5.95%	0.326%
82	12.286	1731	1734	1737	rBV2	108832	106176	8.95%	0.490%
83	12.339	1741	1743	1746	rVV	126933	113060	9.53%	0.522%
84	12.368	1746	1748	1752	rVB2	79455	81886	6.90%	0.378%
85	12.445	1758	1761	1764	rBV	523662	430833	36.31%	1.987%
86	12.557	1778	1780	1786	rBV5	50410	64186	5.41%	0.296%
87	12.668	1796	1799	1802	rVB	365331	323479	27.26%	1.492%
88	12.710	1805	1806	1809	rVB2	56237	43019	3.63%	0.198%
89	12.815	1821	1824	1828	rVB	763145	739435	62.32%	3.411%
90	13.104	1870	1873	1875	rBV	105045	99867	8.42%	0.461%
91	13.410	1923	1925	1930	rVB2	103051	100244	8.45%	0.462%
92	13.504	1939	1941	1944	rBV	108295	83784	7.06%	0.386%
93	13.680	1969	1971	1974	rVB	118254	100906	8.50%	0.465%
94	13.868	1998	2003	2006	rBV2	822546	1128476	95.10%	5.206%
95	13.892	2006	2007	2009	rVB	295765	162683	13.71%	0.750%
96	14.009	2025	2027	2031	rVB	107590	80920	6.82%	0.373%
97	14.051	2031	2034	2036	rBV2	226195	195096	16.44%	0.900%
98	14.886	2173	2176	2178	rBV	241548	290541	24.49%	1.340%
99	15.162	2221	2223	2229	rVB	245650	306326	25.82%	1.413%
100	15.286	2241	2244	2247	rVB	312787	319620	26.94%	1.474%

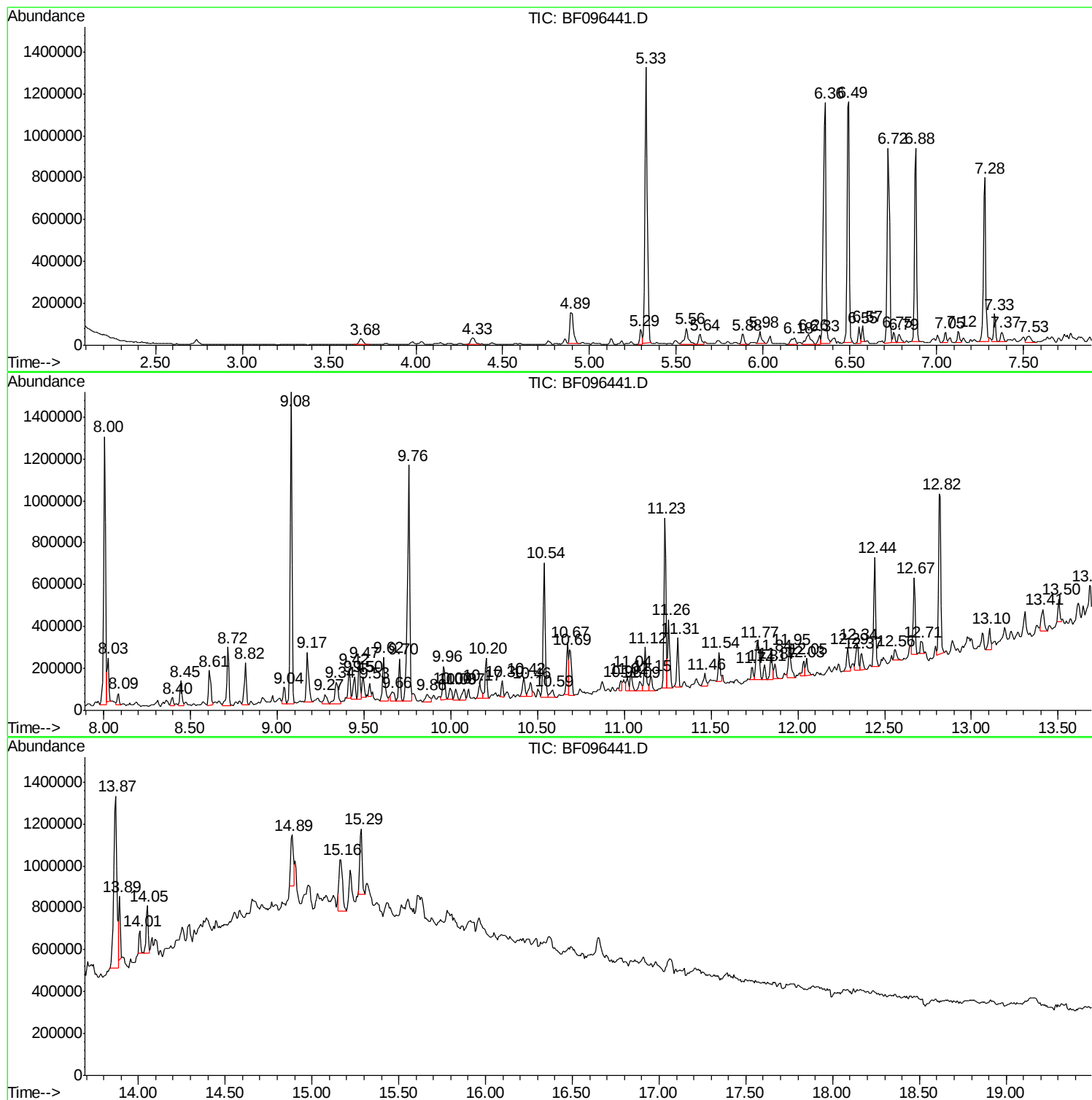
Sum of corrected areas: 21677637

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Instrument :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF070117.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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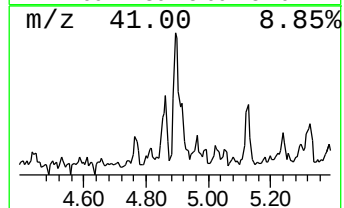
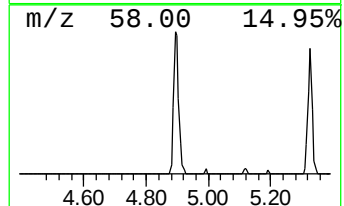
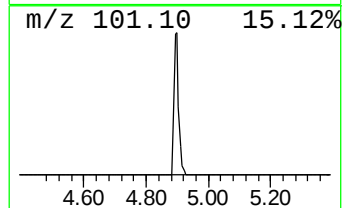
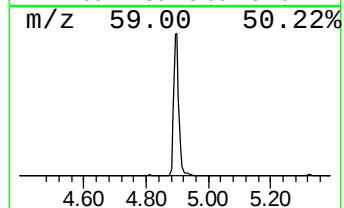
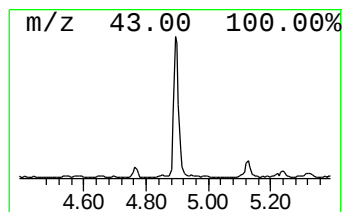
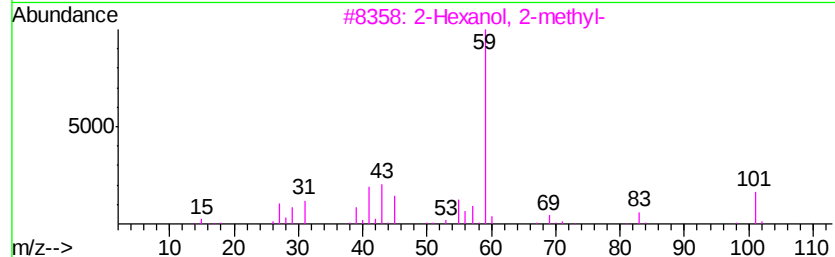
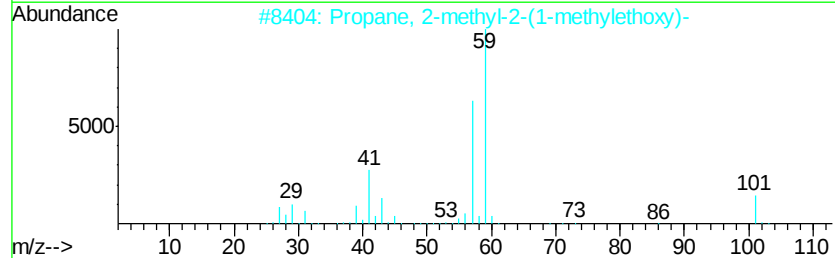
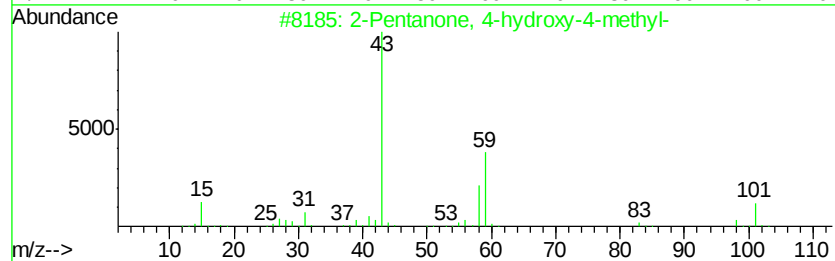
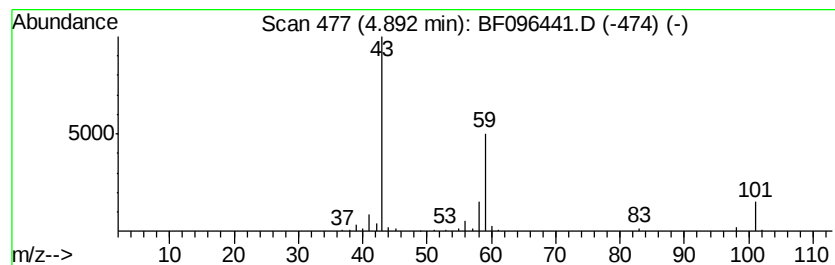
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF070117.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.89	5.05 ng	188595	1,4-Dichlorobenzene-d4	6.72

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	53	
2	Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	36	
3	2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	28	
4	1,2-Butanediol	90	C4H10O2	000584-03-2	28	
5	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	9	



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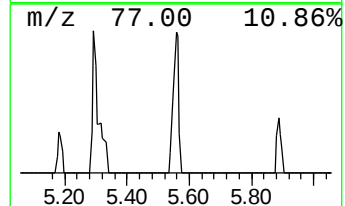
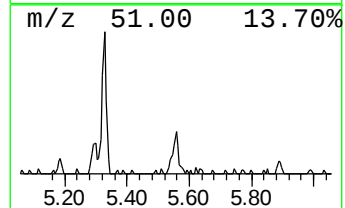
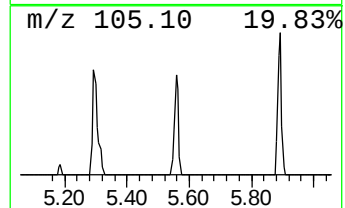
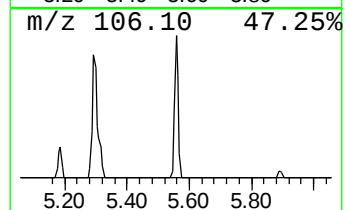
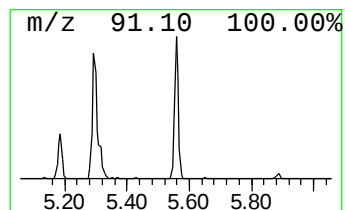
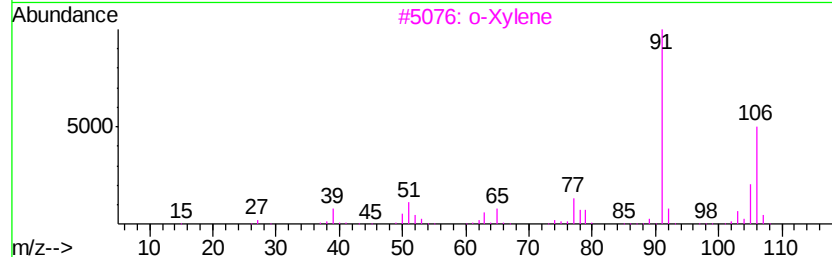
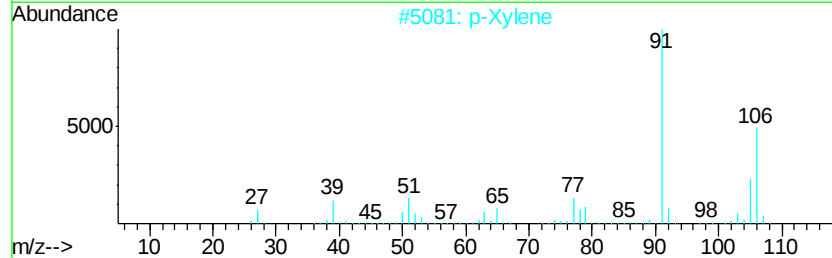
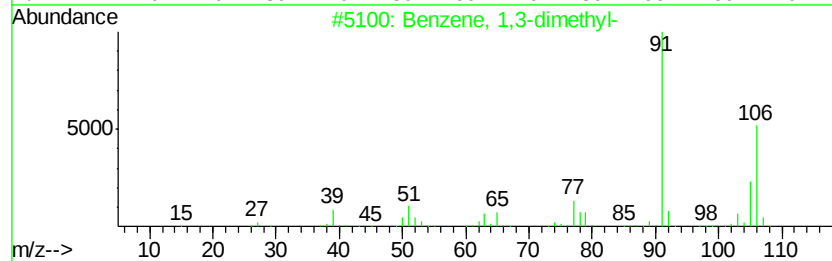
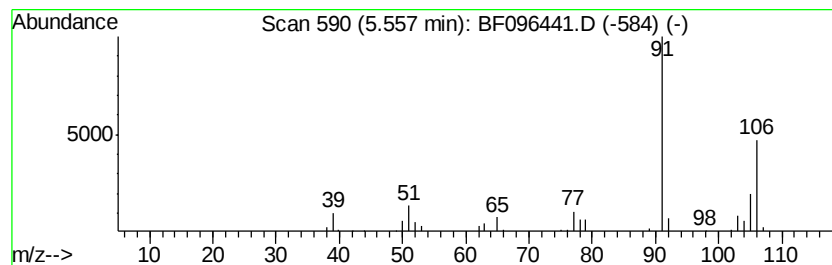
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 Benzene, 1,3-dimethyl- Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.56	3.05 ng	113923	1,4-Dichlorobenzene-d4	6.72

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	94
2			p-Xylene	106	C8H10	000106-42-3	94
3			o-Xylene	106	C8H10	000095-47-6	94
4			1,3-Cyclopentadiene, 5-(1-methyl...	106	C8H10	002175-91-9	83
5			Ethylbenzene	106	C8H10	000100-41-4	76



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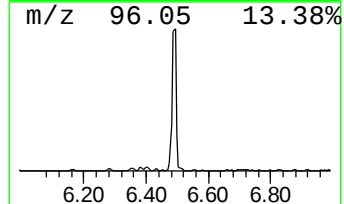
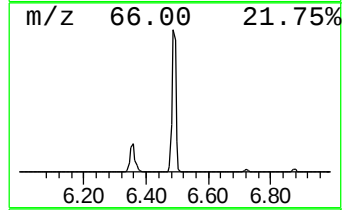
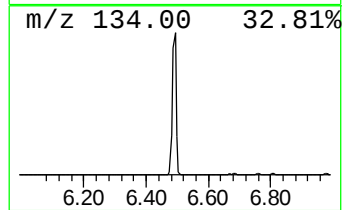
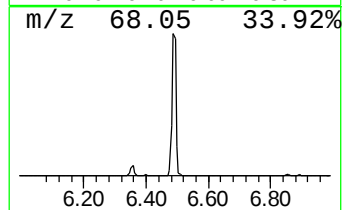
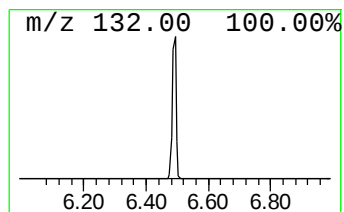
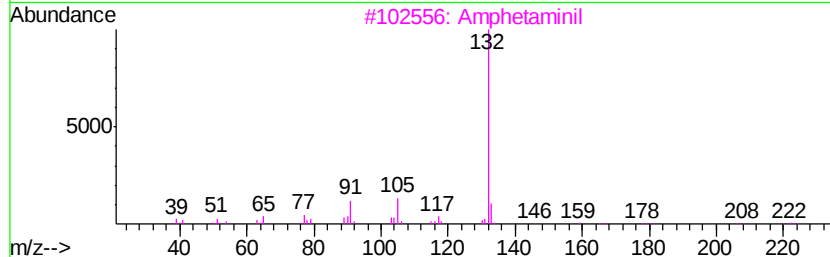
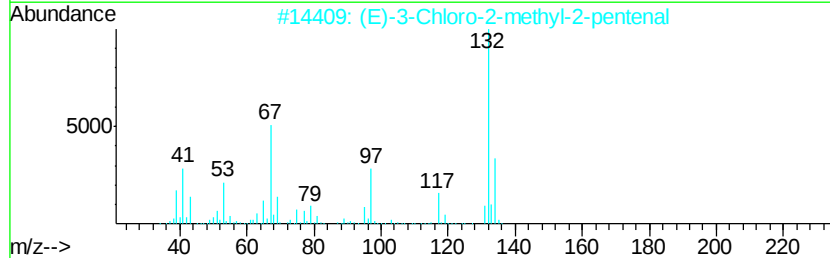
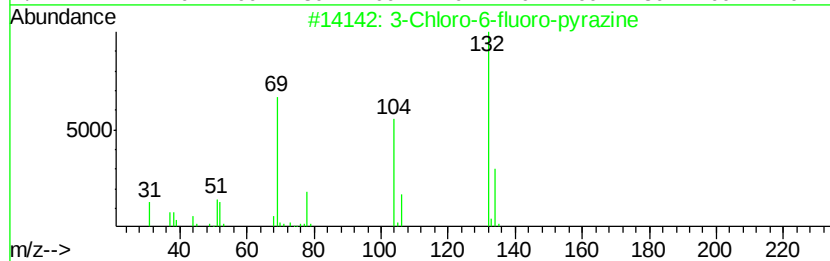
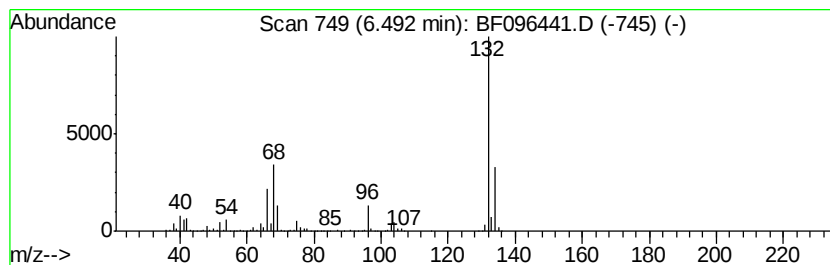
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 unknown6.49 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.49	28.45 ng	1062030	1,4-Dichlorobenzene-d4	6.72

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	38
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	16
3		Amphetaminil	250	C17H18N2	017590-01-1	12
4		1,3,5-Trifluorobenzene	132	C6H3F3	000372-38-3	9
5		Xenon	132	Xe	007440-63-3	9



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Operator : SJ/MA
Sample : I3930-08 2X
Misc :
ALS Vial : 12 Sample Multiplier: 1

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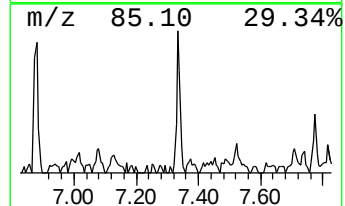
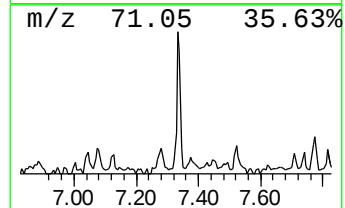
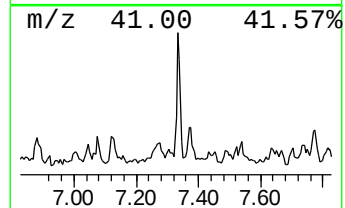
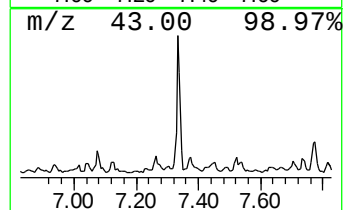
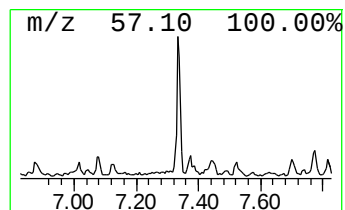
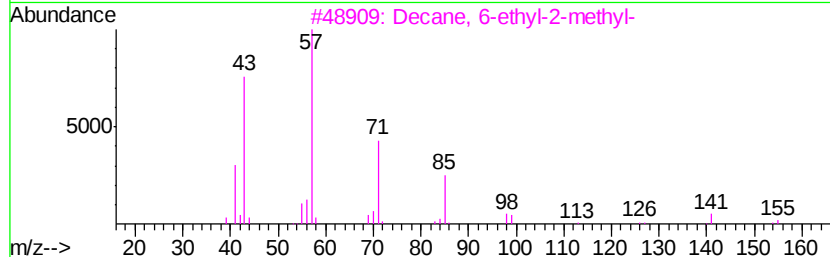
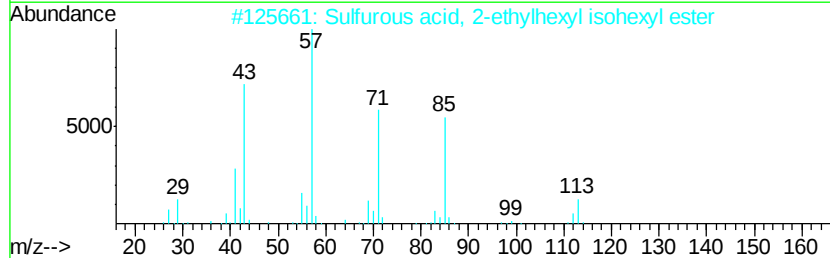
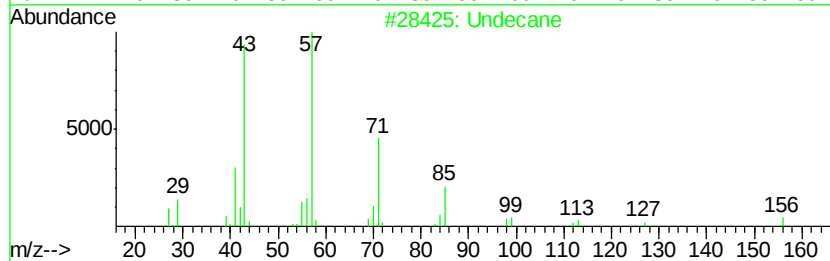
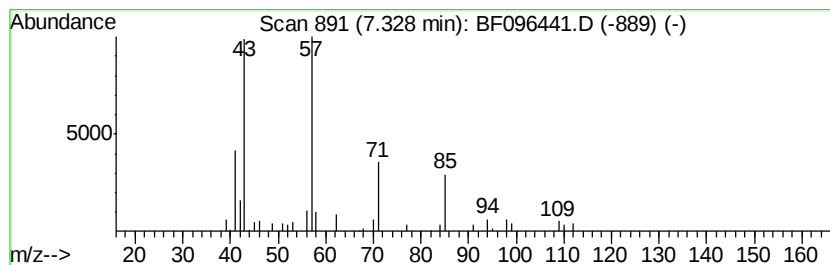
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Undecane Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.33	3.01 ng	112328	1,4-Dichlorobenzene-d4	6.72

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Undecane		156	C11H24	001120-21-4	87
2	Sulfurous acid, 2-ethylhexyl iso...		278	C14H30O3S	1000309-19-0	78
3	Decane, 6-ethyl-2-methyl-		184	C13H28	062108-21-8	78
4	Tetradecane		198	C14H30	000629-59-4	78
5	Sulfurous acid, hexyl octyl ester		278	C14H30O3S	1000309-13-0	64



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF070317\
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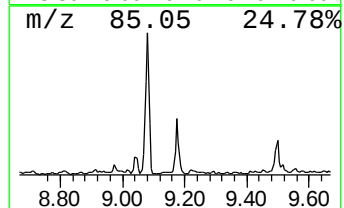
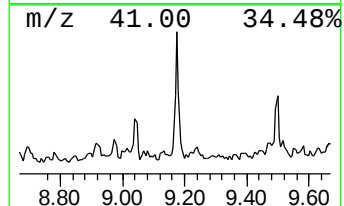
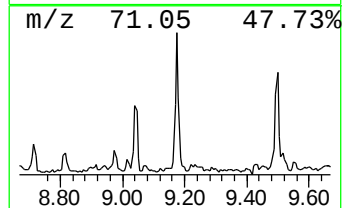
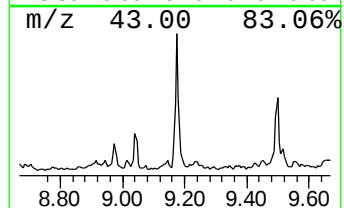
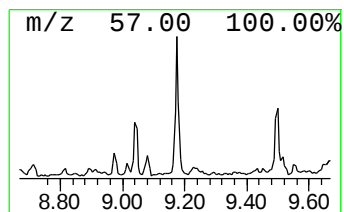
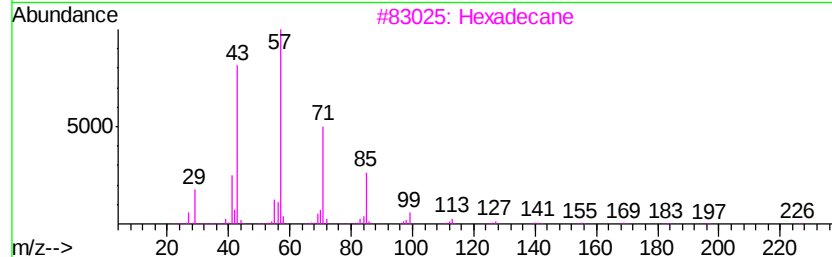
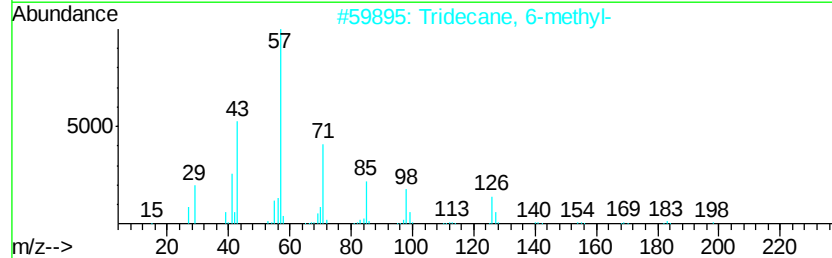
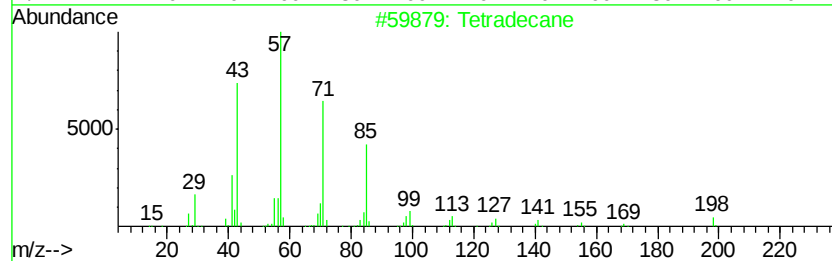
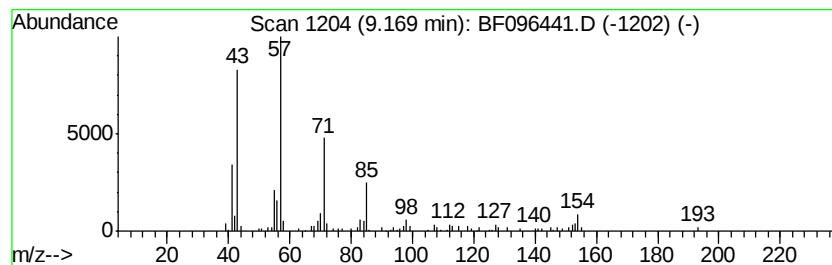
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Tetradecane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.17	4.22 ng	208359	Acenaphthene-d10	9.76
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Tetradecane	198 C14H30	000629-59-4 70
2		Tridecane, 6-methyl-	198 C14H30	013287-21-3 64
3		Hexadecane	226 C16H34	000544-76-3 62
4		Tridecane	184 C13H28	000629-50-5 58
5		Pentadecane	212 C15H32	000629-62-9 58



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF070317\
Data File : BF096441.D
Acq On : 3 Jul 2017 15:14
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ClientSampleId :
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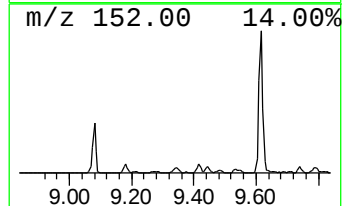
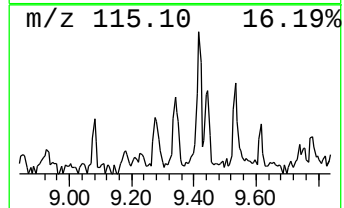
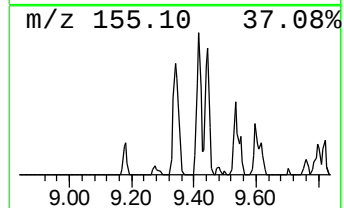
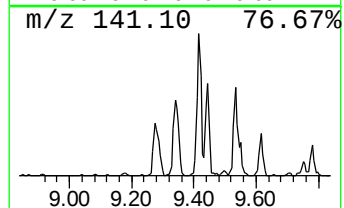
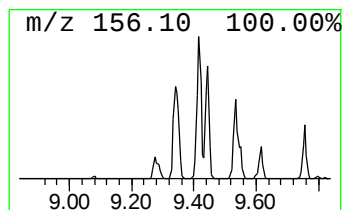
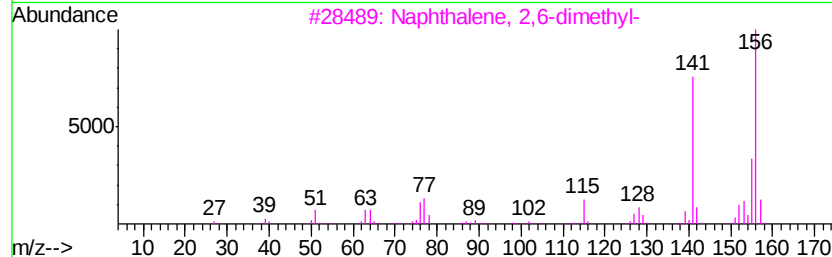
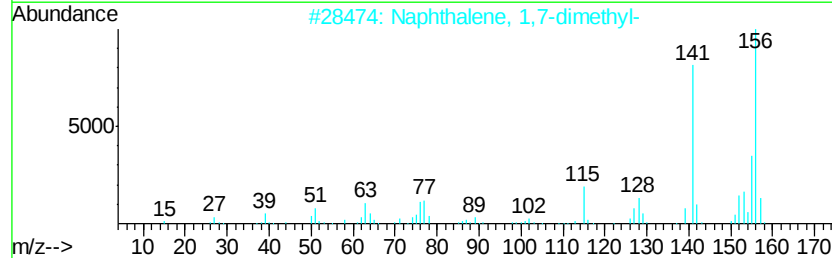
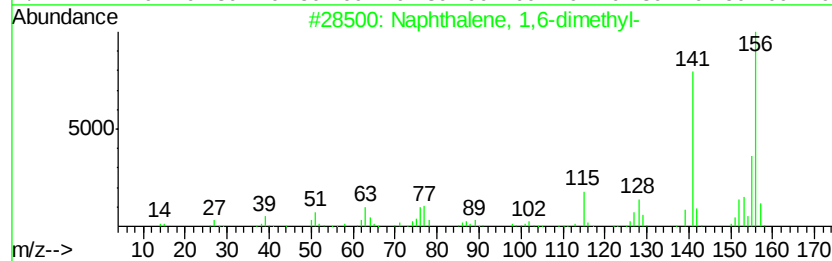
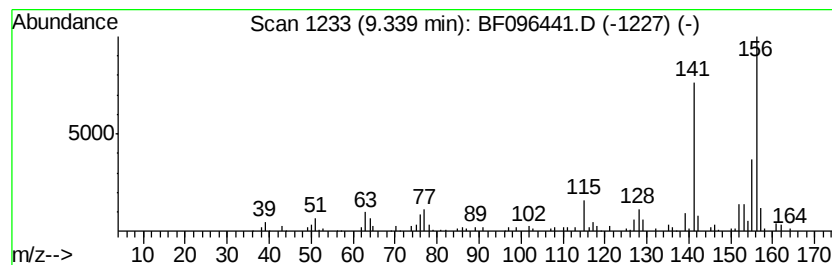
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Naphthalene, 1,6-dimethyl- Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.34	3.05 ng	150628	Acenaphthene-d10	9.76

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	98
3			Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	96
4			Naphthalene, 1,8-dimethyl-	156	C12H12	000569-41-5	96
5			Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	96



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF070317\
Data File : BF096441.D
Acq On : 3 Jul 2017 15:14
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Misc :
ALS Vial : 12 Sample Multiplier: 1

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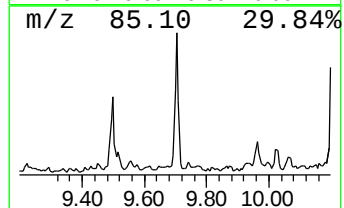
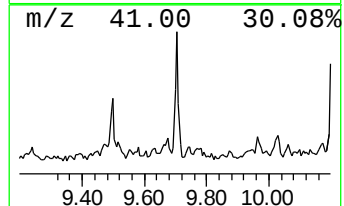
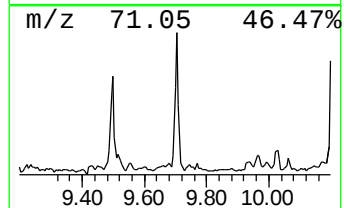
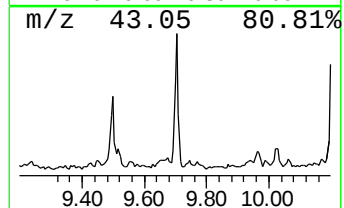
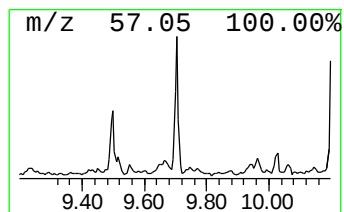
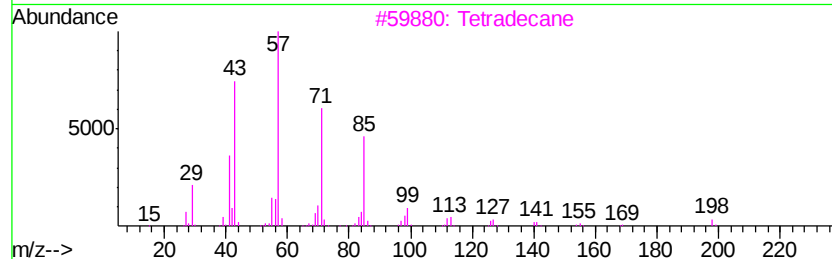
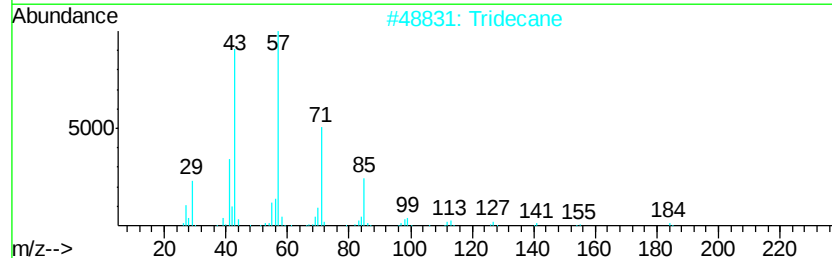
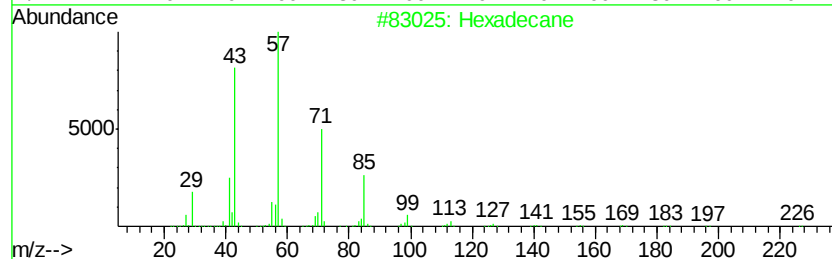
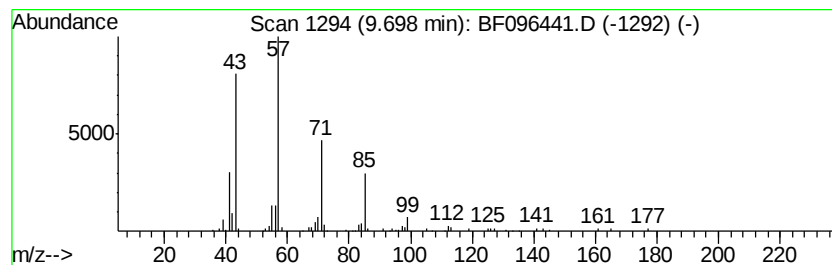
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Hexadecane Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.70	3.21 ng	158601	Acenaphthene-d10	9.76

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecane	226	C16H34	000544-76-3	91
2		Tridecane	184	C13H28	000629-50-5	90
3		Tetradecane	198	C14H30	000629-59-4	80
4		Undecane	156	C11H24	001120-21-4	80
5		Dodecane, 2-methyl-	184	C13H28	001560-97-0	72



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF070317\
Data File : BF096441.D
Acq On : 3 Jul 2017 15:14
Operator : SJ/MA
Sample : I3930-08 2X
Misc :
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Instrument :
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ClientSampleId :
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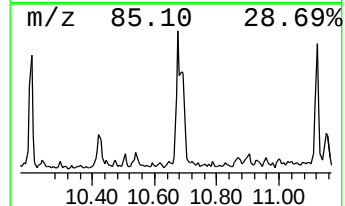
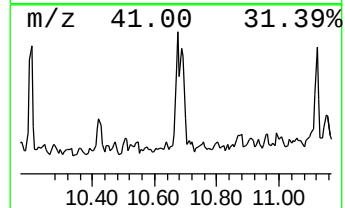
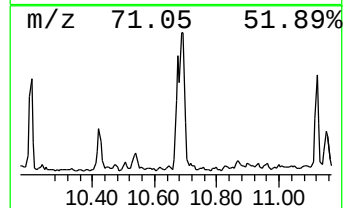
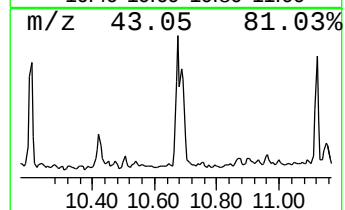
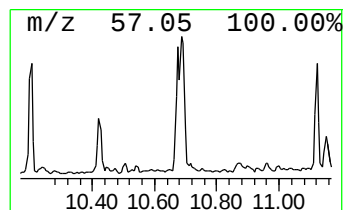
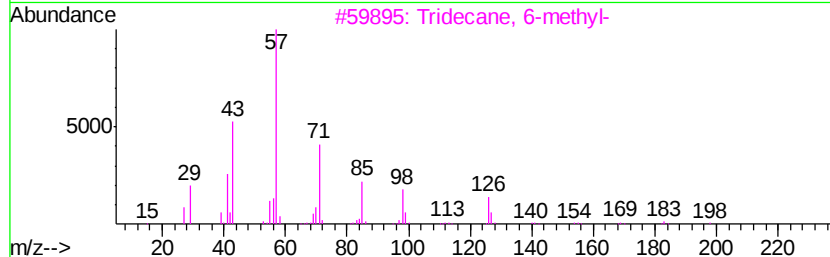
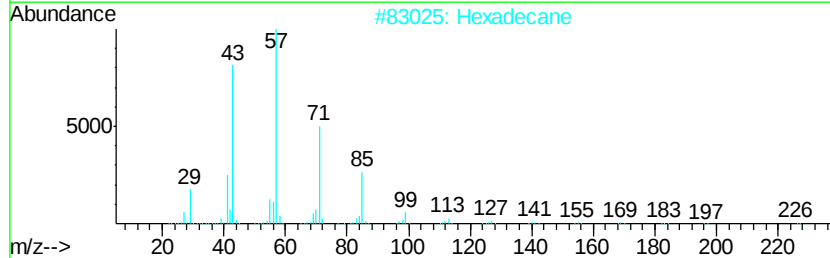
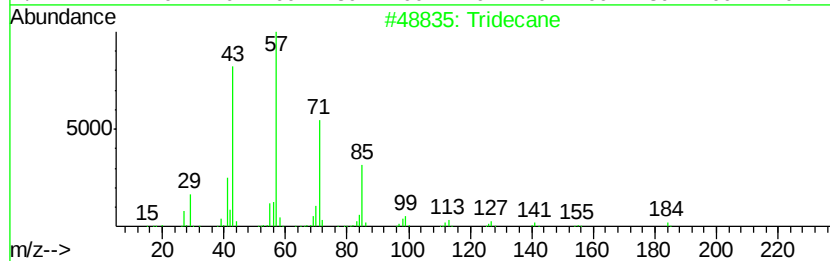
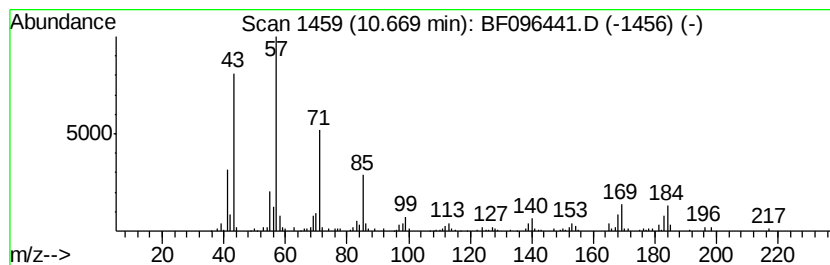
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 Tridecane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.67	6.67 ng	228531	Phenanthrene-d10	11.23

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tridecane	184	C13H28	000629-50-5	86
2			Hexadecane	226	C16H34	000544-76-3	83
3			Tridecane, 6-methyl-	198	C14H30	013287-21-3	72
4			Undecane, 3-methyl-	170	C12H26	001002-43-3	64
5			Tetradecane	198	C14H30	000629-59-4	64



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF070317\
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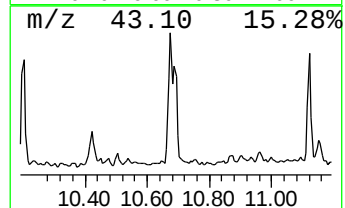
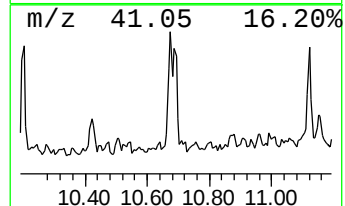
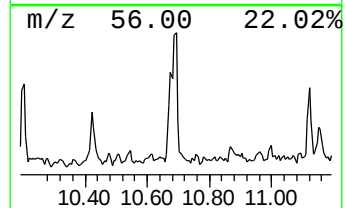
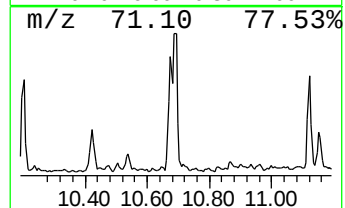
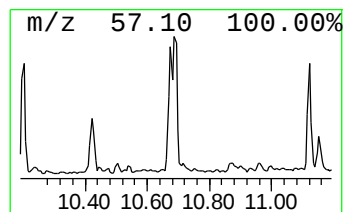
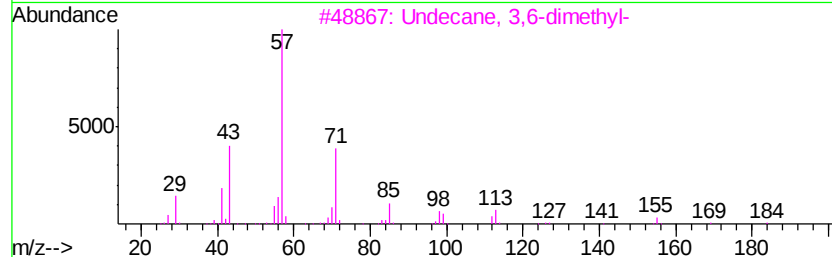
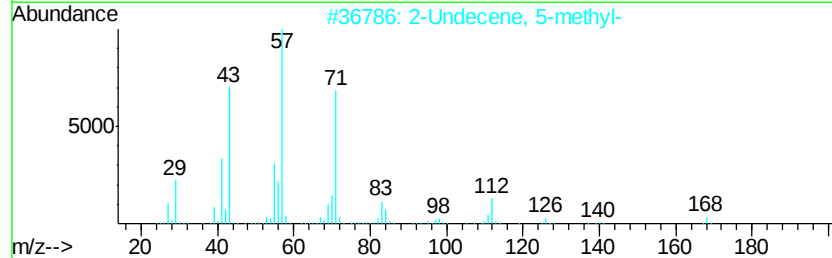
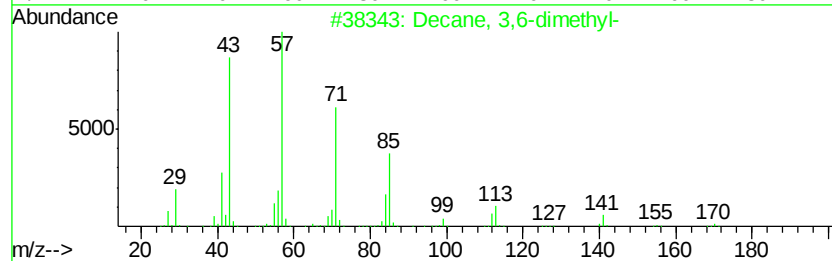
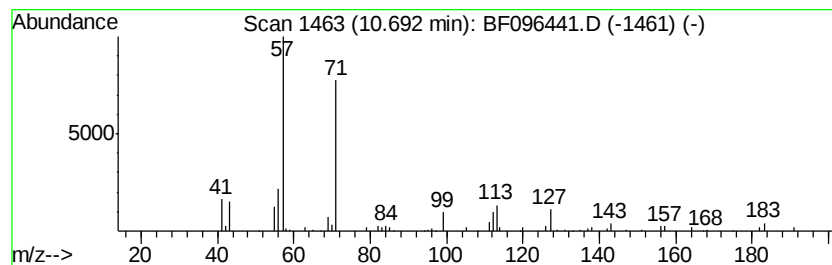
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Decane, 3,6-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD		R.T.
10.69	5.20 ng	178327	Phenanthrene-d10		11.23
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Decane, 3,6-dimethyl-	170	C12H26	017312-53-7	64
2	2-Undecene, 5-methyl-	168	C12H24	056851-34-4	53
3	Undecane, 3,6-dimethyl-	184	C13H28	017301-28-9	53
4	Undecane, 6-ethyl-	184	C13H28	017312-60-6	53
5	Undecane, 4,6-dimethyl-	184	C13H28	017312-82-2	50



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF070317\
Data File : BF096441.D
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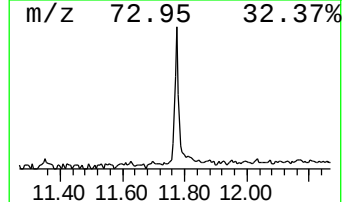
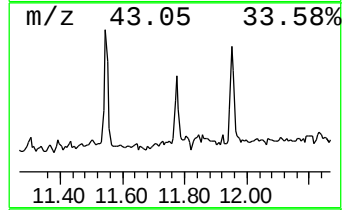
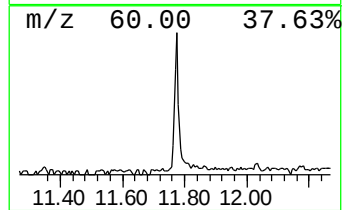
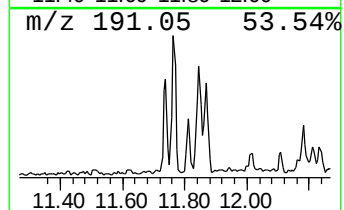
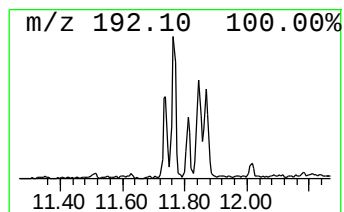
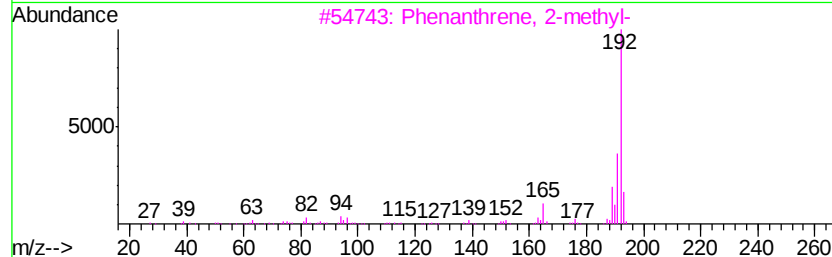
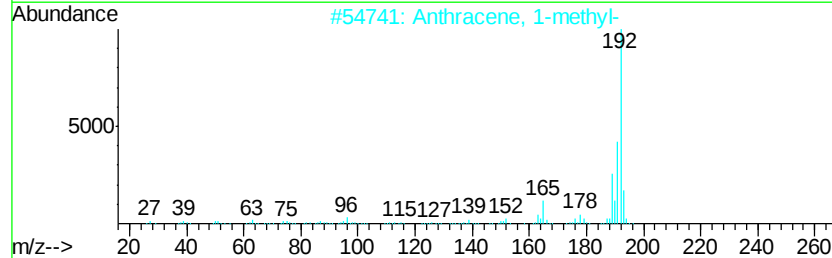
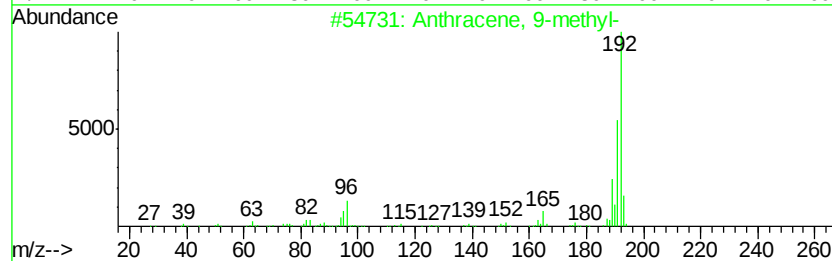
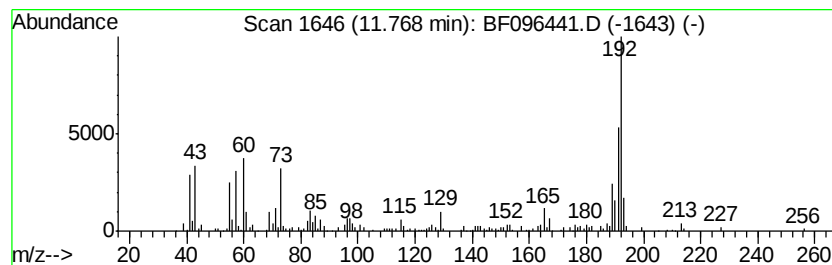
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Anthracene, 9-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.77	6.30 ng	216023	Phenanthrene-d10	11.23

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Anthracene, 9-methyl-	192	C15H12	000779-02-2	83
2		Anthracene, 1-methyl-	192	C15H12	000610-48-0	83
3		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	83
4		Anthracene, 2-methyl-	192	C15H12	000613-12-7	83
5		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	78



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF070317\
Data File : BF096441.D
Acq On : 3 Jul 2017 15:14
Operator : SJ/MA
Sample : I3930-08 2X
Misc :
ALS Vial : 12 Sample Multiplier: 1

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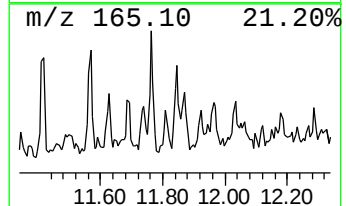
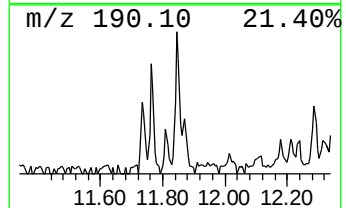
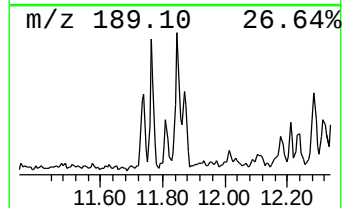
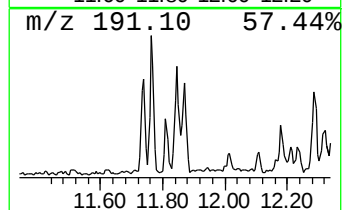
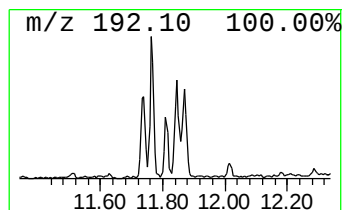
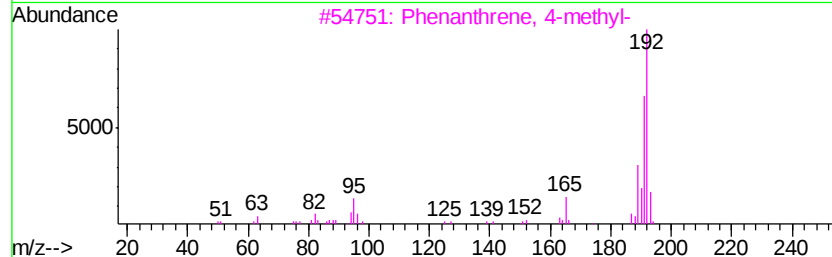
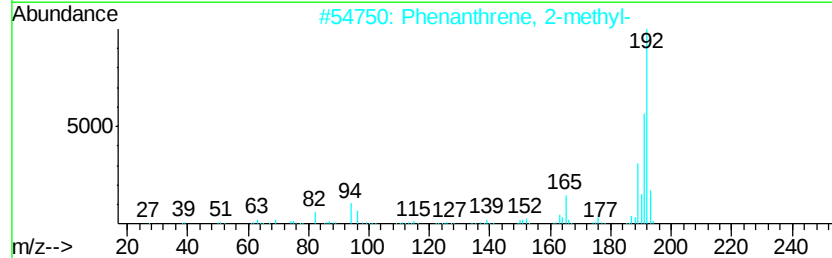
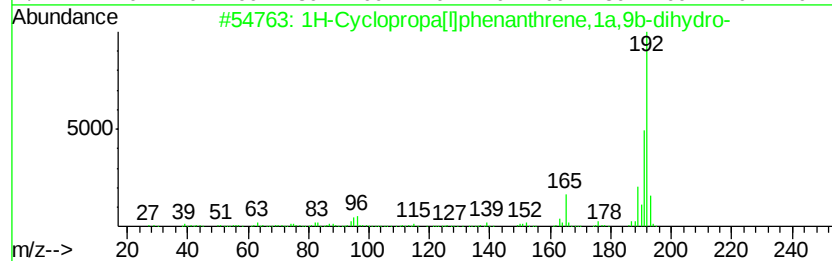
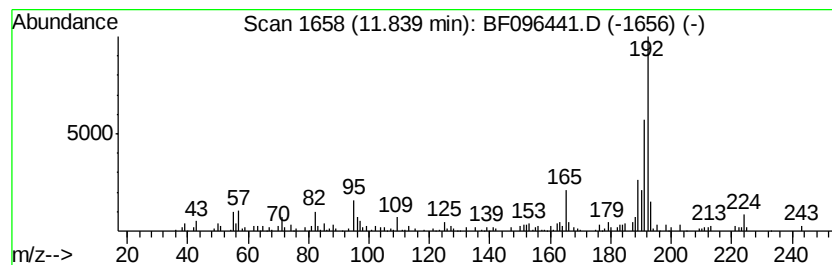
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 1H-Cyclopropa[l]phenanthren... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.84	3.44 ng	117874	Phenanthrene-d10	11.23

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Cyclopropa[l]phenanthrene, 1a, ...	192	C15H12	000949-41-7	93
2			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	92
3			Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	92
4			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	91
5			1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	91



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF070317\
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Sample : I3930-08 2X
Misc :
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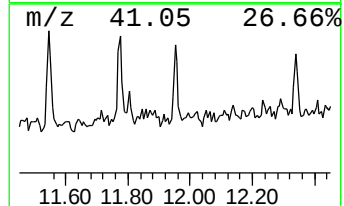
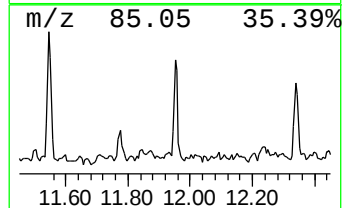
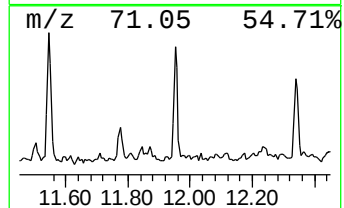
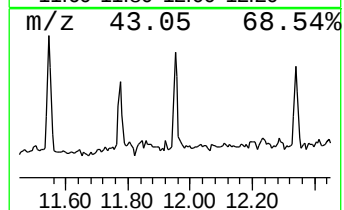
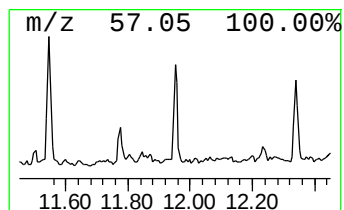
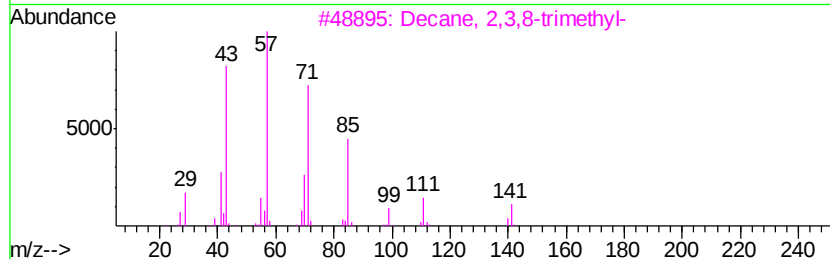
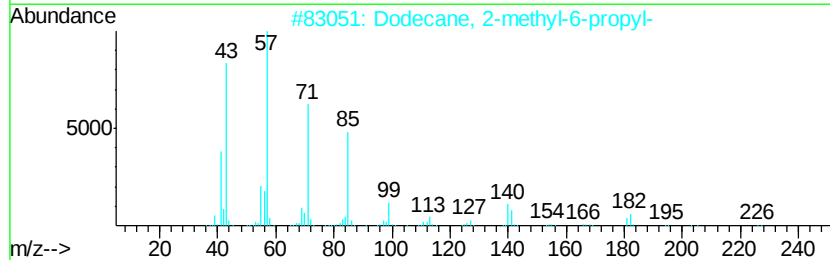
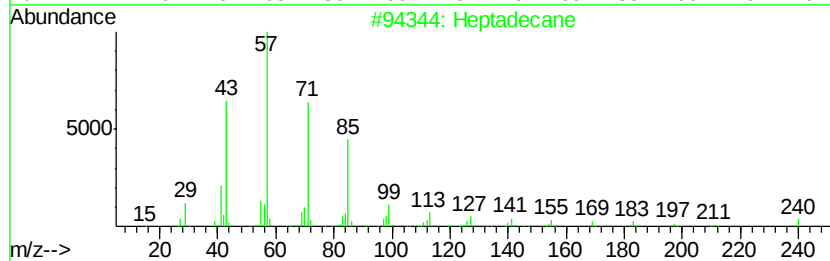
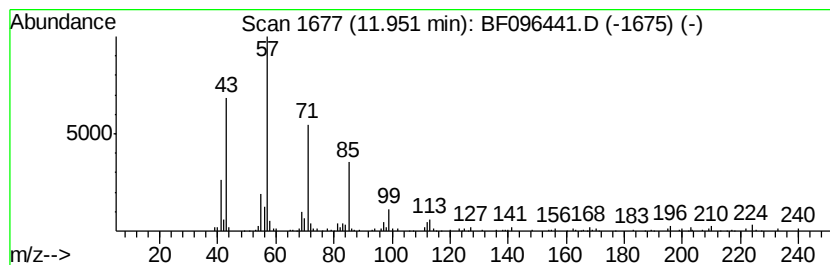
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF070117.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 Heptadecane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD		R.T.	
11.95	4.09 ng	140068	Phenanthrene-d10		11.23	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptadecane	240	C17H36	000629-78-7	80
2		Dodecane, 2-methyl-6-propyl-	226	C16H34	055045-08-4	80
3		Decane, 2,3,8-trimethyl-	184	C13H28	062238-14-6	72
4		Dodecane, 3-methyl-	184	C13H28	017312-57-1	72
5		Octane, 5-ethyl-2-methyl-	156	C11H24	062016-18-6	64



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF070317\
Data File : BF096441.D
Acq On : 3 Jul 2017 15:14
Operator : SJ/MA
Sample : I3930-08 2X
Misc :
ALS Vial : 12 Sample Multiplier: 1

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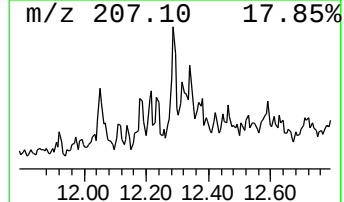
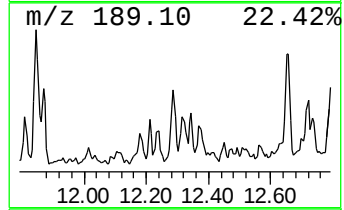
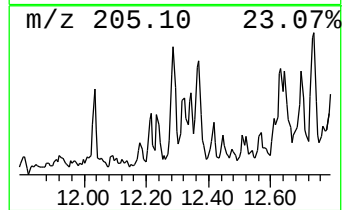
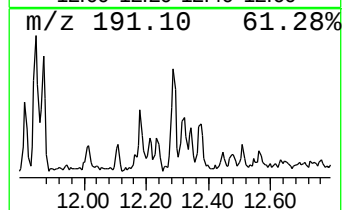
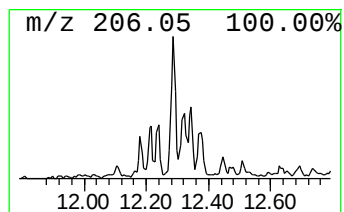
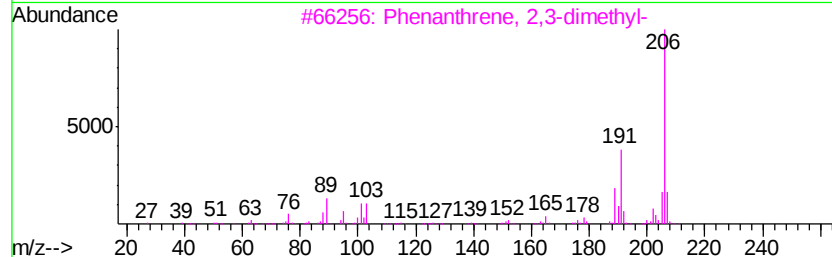
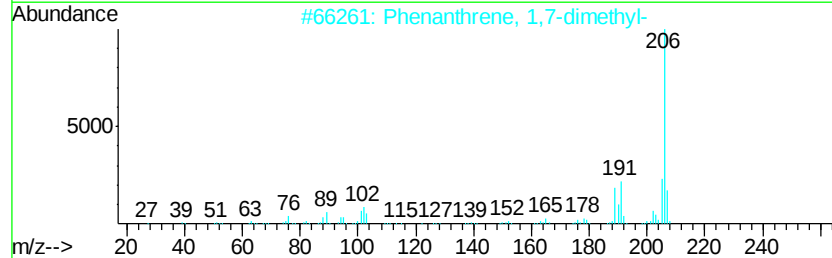
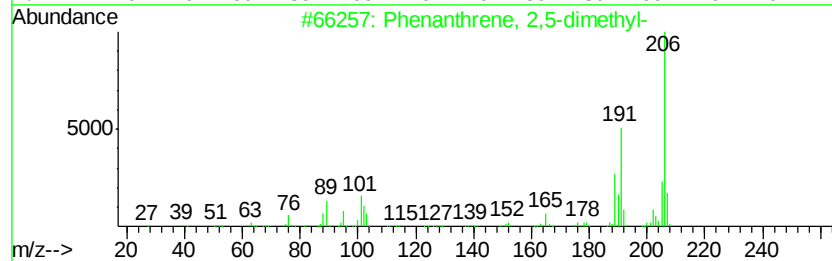
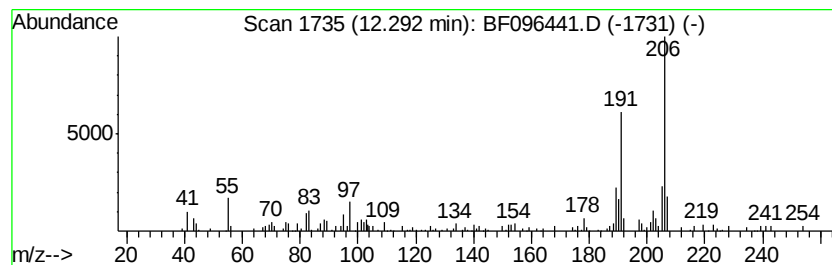
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 Phenanthrene, 2,5-dimethyl- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.29	3.10 ng	106176	Phenanthrene-d10	11.23

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	95
2		Phenanthrene, 1,7-dimethyl-	206	C16H14	000483-87-4	93
3		Phenanthrene, 2,3-dimethyl-	206	C16H14	003674-65-5	93
4		Phenanthrene, 3,6-dimethyl-	206	C16H14	001576-67-6	89
5		1,3-Diphenyl-3-methylcyclopropene	206	C16H14	086544-79-8	89



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF070317\
Data File : BF096441.D
Acq On : 3 Jul 2017 15:14
Operator : SJ/MA
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Misc :
ALS Vial : 12 Sample Multiplier: 1

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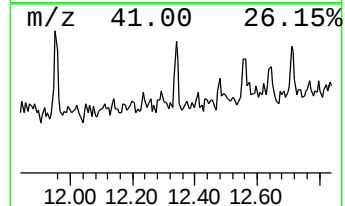
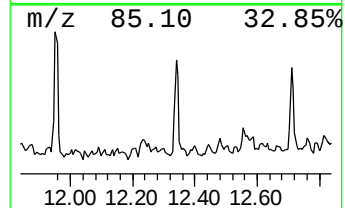
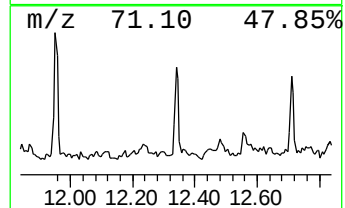
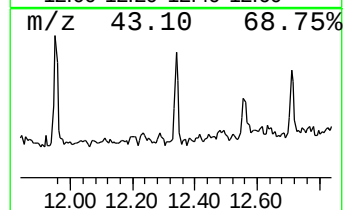
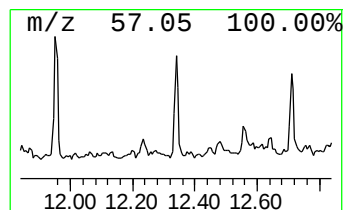
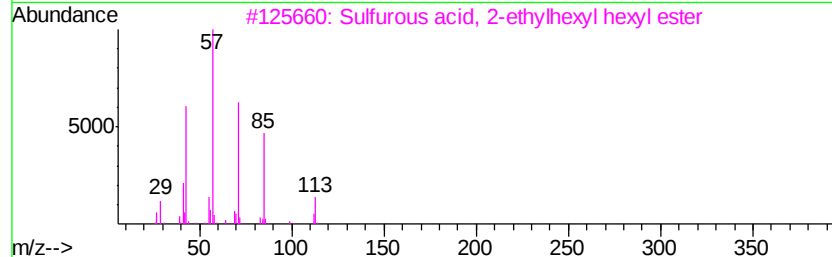
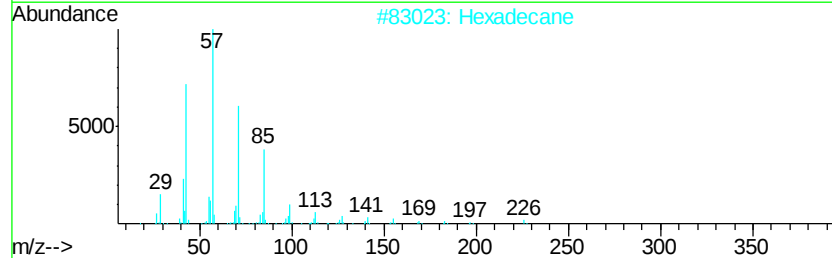
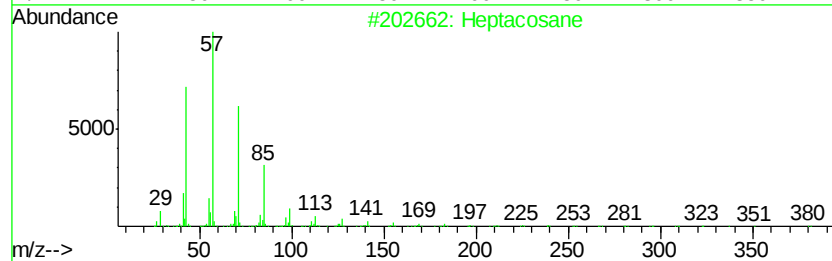
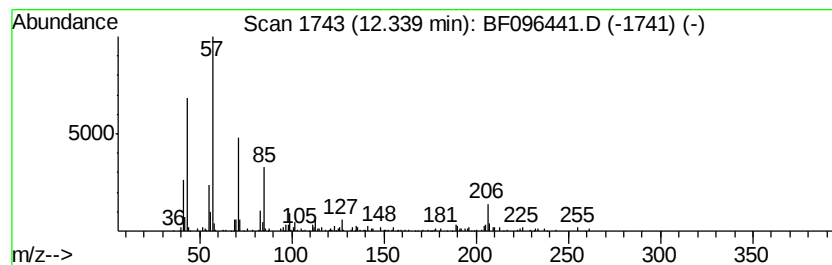
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 Heptacosane Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.34	3.30 ng	113060	Phenanthrene-d10	11.23

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptacosane	380	C27H56	000593-49-7	68
2		Hexadecane	226	C16H34	000544-76-3	64
3		Sulfurous acid, 2-ethylhexyl hex...	278	C14H30O3S	1000309-20-2	59
4		Eicosane	282	C20H42	000112-95-8	52
5		Tetratetracontane	619	C44H90	007098-22-8	52



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF070317\
Data File : BF096441.D
Acq On : 3 Jul 2017 15:14
Operator : SJ/MA
Sample : I3930-08 2X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
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ClientSampleId :
B3-0-2

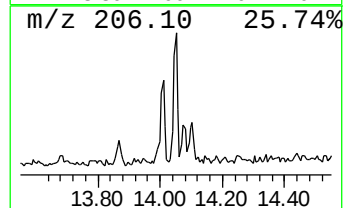
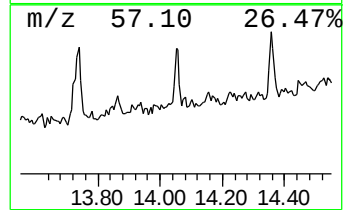
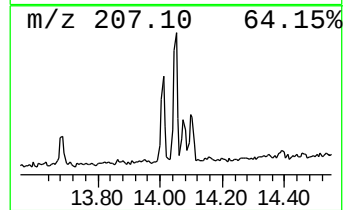
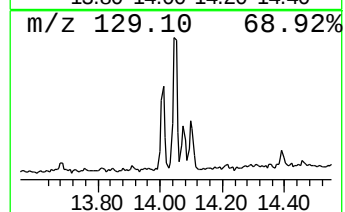
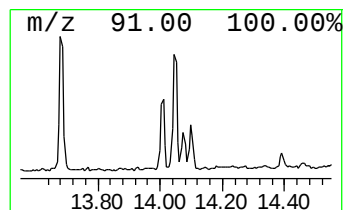
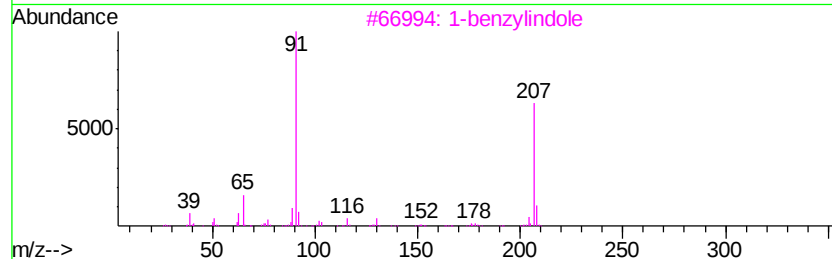
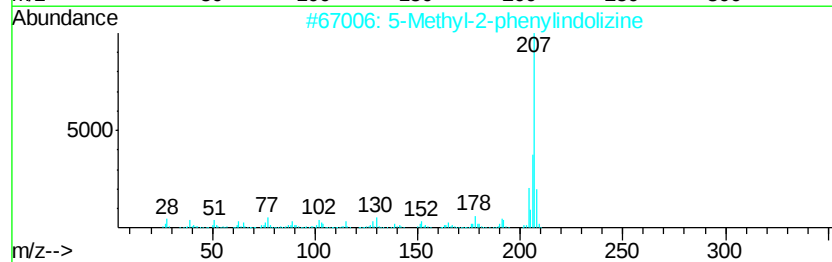
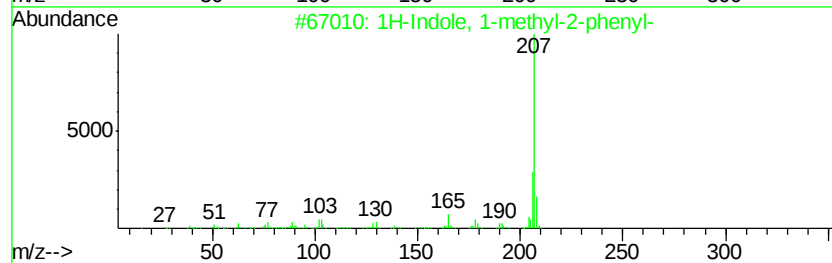
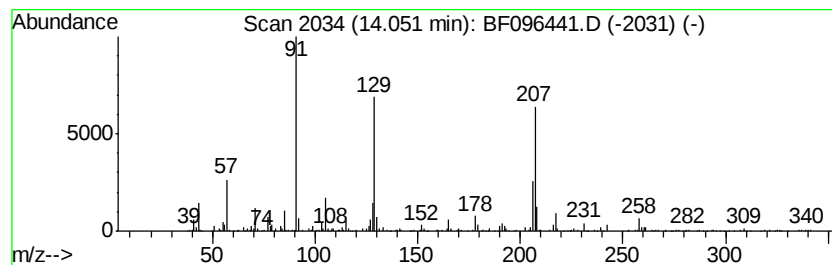
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 19 unknown14.05 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.05	3.46 ng	195096	Chrysene-d12	13.87

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Indole, 1-methyl-2-phenyl-	207	C15H13N	003558-24-5	38
2			5-Methyl-2-phenylindolizine	207	C15H13N	036944-99-7	30
3			1-benzylindole	207	C15H13N	1000334-31-3	27
4			1H-Indole, 2-methyl-3-phenyl-	207	C15H13N	004757-69-1	27
5			(2,3-Diphenylcyclopropyl)methyl ...	332	C22H20OS	131758-71-9	25



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF070317\
Data File : BF096441.D
Acq On : 3 Jul 2017 15:14
Operator : SJ/MA
Sample : I3930-08 2X
Misc :
ALS Vial : 12 Sample Multiplier: 1

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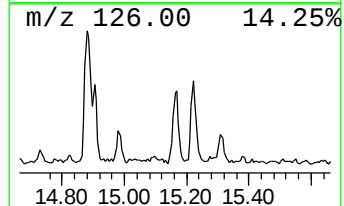
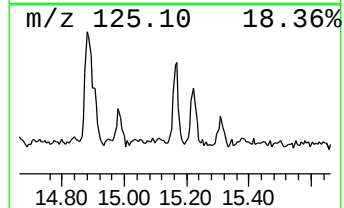
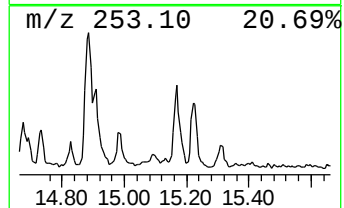
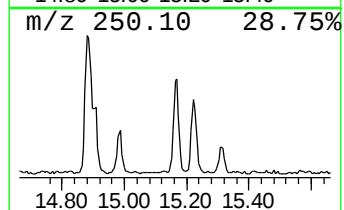
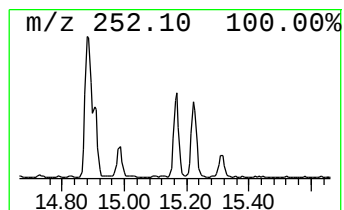
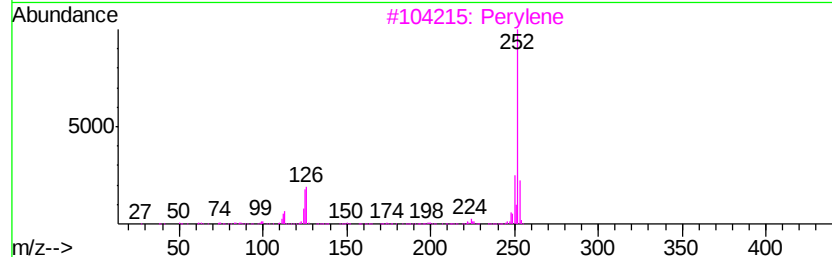
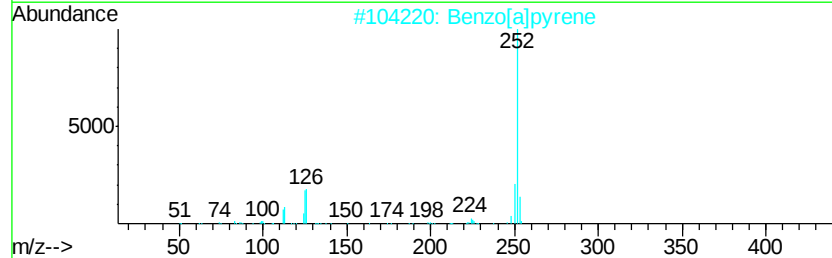
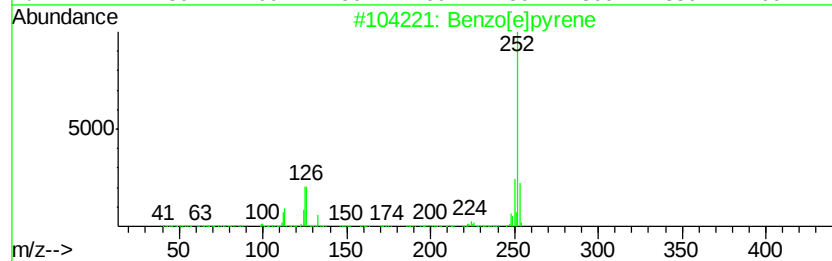
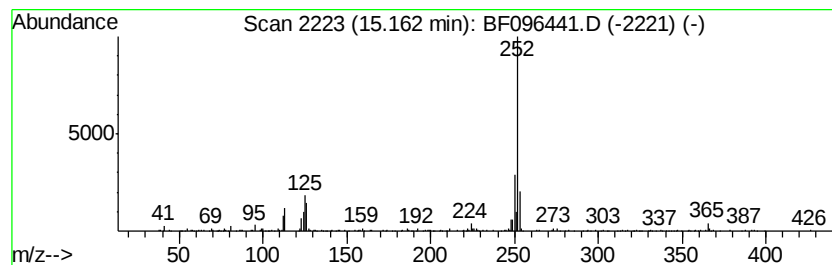
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 20 Benzo[e]pyrene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.16	19.17 ng	306326	Perylene-d12	15.29

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzo[e]pyrene	252	C20H12	000192-97-2	93
2		Benzo[a]pyrene	252	C20H12	000050-32-8	86
3		Perylene	252	C20H12	000198-55-0	76
4		Benzo[j]fluoranthene	252	C20H12	000205-82-3	76
5		Benz[e]acephenanthrylene	252	C20H12	000205-99-2	74



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF070317\
 Data File : BF096441.D
 Acq On : 3 Jul 2017 15:14
 Operator : SJ/MA
 Sample : I3930-08 2X
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 B3-0-2

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF070117.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
2-Pentanone, 4-hy...	4.89	5.0 ng		188595	1	6.72	746604	20.0
Benzene, 1,3-dime...	5.56	3.0 ng		113923	1	6.72	746604	20.0
unknown6.49	6.49	28.4 ng		1062030	1	6.72	746604	20.0
Undecane	7.33	3.0 ng		112328	1	6.72	746604	20.0
Tetradecane	9.17	4.2 ng		208359	3	9.76	988033	20.0
Naphthalene, 1,6-...	9.34	3.0 ng		150628	3	9.76	988033	20.0
Hexadecane	9.70	3.2 ng		158601	3	9.76	988033	20.0
Tridecane	10.67	6.7 ng		228531	4	11.23	685696	20.0
Decane, 3,6-dimet...	10.69	5.2 ng		178327	4	11.23	685696	20.0
Anthracene, 9-met...	11.77	6.3 ng		216023	4	11.23	685696	20.0
1H-Cyclopropa[l]p...	11.84	3.4 ng		117874	4	11.23	685696	20.0
Heptadecane	11.95	4.1 ng		140068	4	11.23	685696	20.0
Phenanthrene, 2,5...	12.29	3.1 ng		106176	4	11.23	685696	20.0
Heptacosane	12.34	3.3 ng		113060	4	11.23	685696	20.0
unknown14.05	14.05	3.5 ng		195096	5	13.87	1128480	20.0
Benzo[e]pyrene	15.16	19.2 ng		306326	6	15.29	319620	20.0