

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF100717\
 Data File : BF099196.D
 Acq On : 7 Oct 2017 16:07
 Operator : SJ/JU
 Sample : I5586-19 2X
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 G56A-3-6

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-BF092317.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	2.122	3	6	13	rVB	70768	90343	6.34%	0.383%
2	3.004	144	156	162	rBV2	28678	54949	3.86%	0.233%
3	3.934	310	314	321	rVB	52446	67400	4.73%	0.286%
4	5.075	505	508	515	rBV	215344	243427	17.09%	1.033%
5	5.451	568	572	574	rBV	96511	89692	6.30%	0.380%
6	5.481	574	577	587	rVB	1276376	1211986	85.10%	5.141%
7	5.710	613	616	620	rBV	89181	84589	5.94%	0.359%
8	6.022	664	669	673	rVB3	34448	43192	3.03%	0.183%
9	6.116	678	685	691	rBV7	33894	66732	4.69%	0.283%
10	6.386	728	731	734	rBV2	57012	56750	3.98%	0.241%
11	6.492	746	749	754	rVV	1280987	1205741	84.66%	5.114%
12	6.633	769	773	776	rBV	1562403	1274783	89.51%	5.407%
13	6.686	779	782	786	rBV2	116885	113886	8.00%	0.483%
14	6.863	809	812	816	rVV	952674	790041	55.47%	3.351%
15	6.922	818	822	825	rBV	61593	61461	4.32%	0.261%
16	7.022	833	839	842	rBV	1159619	1048931	73.65%	4.449%
17	7.133	852	858	861	rVB3	35175	43892	3.08%	0.186%
18	7.257	873	879	884	rBV4	36099	61397	4.31%	0.260%
19	7.428	904	908	910	rVV	822563	744252	52.26%	3.157%
20	7.451	910	912	916	rVB	90711	72404	5.08%	0.307%
21	7.504	916	921	924	rBV5	42177	49106	3.45%	0.208%
22	8.122	1022	1026	1027	rBV	87916	77068	5.41%	0.327%
23	8.145	1027	1030	1032	rVV	1209370	981512	68.91%	4.163%
24	8.169	1032	1034	1036	rVV	261860	196440	13.79%	0.833%
25	8.198	1036	1039	1041	rVB	70085	55491	3.90%	0.235%
26	8.433	1072	1079	1083	rBV4	34271	45886	3.22%	0.195%
27	8.557	1097	1100	1103	rVV	119302	101785	7.15%	0.432%
28	8.727	1126	1129	1132	rBV	117607	83111	5.84%	0.353%
29	8.857	1148	1151	1154	rBV	319569	235343	16.52%	0.998%
30	8.957	1164	1168	1171	rBV	224272	171098	12.01%	0.726%
31	9.157	1200	1202	1204	rVB	66782	46138	3.24%	0.196%
32	9.216	1208	1212	1216	rBV	1635992	1424247	100.00%	6.041%
33	9.292	1222	1225	1228	rBV	134683	114026	8.01%	0.484%
34	9.316	1228	1229	1233	rVV	55329	44878	3.15%	0.190%

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Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF092317.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	9.416	1243	1246	1250	rVB	37955	50389	3.54%	0.214%
36	9.486	1253	1258	1263	rVB2	113287	133876	9.40%	0.568%
37	9.557	1267	1270	1272	rBV	154049	150399	10.56%	0.638%
38	9.586	1272	1275	1276	rVV	105679	86822	6.10%	0.368%
39	9.610	1276	1279	1286	rVB	327331	301034	21.14%	1.277%
40	9.674	1286	1290	1292	rBV	104330	88970	6.25%	0.377%
41	9.763	1302	1305	1307	rVB	96605	86235	6.05%	0.366%
42	9.822	1307	1315	1318	rBV	151135	127241	8.93%	0.540%
43	9.880	1321	1325	1326	rVV	71333	74790	5.25%	0.317%
44	9.898	1326	1328	1332	rVV	992896	942542	66.18%	3.998%
45	9.933	1332	1334	1337	rVV3	52999	49470	3.47%	0.210%
46	10.004	1342	1346	1350	rBV	44090	55081	3.87%	0.234%
47	10.051	1350	1354	1357	rBV4	40102	60503	4.25%	0.257%
48	10.104	1359	1363	1366	rVV2	150726	151928	10.67%	0.644%
49	10.139	1366	1369	1373	rVV2	89721	87782	6.16%	0.372%
50	10.216	1379	1382	1384	rBV	75139	68009	4.78%	0.288%
51	10.245	1384	1387	1389	rVV	60412	53758	3.77%	0.228%
52	10.321	1392	1400	1403	rBV2	202332	241385	16.95%	1.024%
53	10.439	1417	1420	1423	rBV2	90330	81235	5.70%	0.345%
54	10.539	1433	1437	1439	rBV	58987	62771	4.41%	0.266%
55	10.604	1443	1448	1452	rBV2	70882	94540	6.64%	0.401%
56	10.686	1459	1462	1467	rVV	702179	609026	42.76%	2.583%
57	10.804	1477	1482	1489	rVB3	226620	365631	25.67%	1.551%
58	10.992	1508	1514	1517	rBV5	31443	61464	4.32%	0.261%
59	11.121	1532	1536	1539	rBV3	54984	82538	5.80%	0.350%
60	11.192	1545	1548	1551	rBV2	95783	84972	5.97%	0.360%
61	11.239	1552	1556	1559	rVV2	166818	168302	11.82%	0.714%
62	11.268	1559	1561	1569	rVB3	59594	99090	6.96%	0.420%
63	11.386	1577	1581	1583	rBV	925561	774742	54.40%	3.286%
64	11.410	1583	1585	1588	rVV	343148	280341	19.68%	1.189%
65	11.463	1591	1594	1596	rVV	105318	84827	5.96%	0.360%
66	11.598	1614	1617	1619	rVV	57374	64117	4.50%	0.272%
67	11.616	1619	1620	1624	rVV2	50993	56084	3.94%	0.238%
68	11.668	1626	1629	1633	rVV2	121194	115969	8.14%	0.492%
69	11.715	1635	1637	1642	rVB3	47175	52183	3.66%	0.221%
70	11.886	1664	1666	1668	rBV	86412	79044	5.55%	0.335%
71	11.915	1668	1671	1674	rVB	164761	151091	10.61%	0.641%

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72	11.998	1682	1685	1687	rBV2	84733	86398	6.07%	0.366%
73	12.021	1687	1689	1692	rVB	74331	58023	4.07%	0.246%
74	12.074	1695	1698	1700	rBV	116564	92431	6.49%	0.392%
75	12.180	1713	1716	1719	rBV	89606	83851	5.89%	0.356%
76	12.210	1719	1721	1727	rVB2	67912	69741	4.90%	0.296%
77	12.363	1744	1747	1749	rVV2	60752	58557	4.11%	0.248%
78	12.439	1757	1760	1762	rVV	95817	111076	7.80%	0.471%
79	12.463	1762	1764	1768	rVV2	125868	136356	9.57%	0.578%
80	12.527	1772	1775	1778	rVV	73918	82134	5.77%	0.348%
81	12.598	1784	1787	1792	rBV	481628	418519	29.39%	1.775%
82	12.645	1792	1795	1797	rBV	48396	51302	3.60%	0.218%
83	12.833	1821	1827	1828	rBV2	437734	489493	34.37%	2.076%
84	12.851	1828	1830	1833	rVB	428866	373346	26.21%	1.584%
85	12.968	1846	1850	1853	rBV	1066459	859527	60.35%	3.646%
86	13.039	1860	1862	1870	rVB7	50769	90123	6.33%	0.382%
87	13.127	1875	1877	1879	rBV2	52124	57828	4.06%	0.245%
88	13.186	1885	1887	1891	rBV2	63018	68128	4.78%	0.289%
89	13.445	1928	1931	1939	rVB	244519	236778	16.62%	1.004%
90	13.533	1943	1946	1949	rBV3	64083	68378	4.80%	0.290%
91	13.662	1966	1968	1972	rVB	88007	86342	6.06%	0.366%
92	13.774	1985	1987	1990	rVB2	58651	48524	3.41%	0.206%
93	13.857	1999	2001	2008	rVB2	115483	153573	10.78%	0.651%
94	13.998	2022	2025	2026	rBV	63602	63308	4.45%	0.269%
95	14.027	2026	2030	2033	rVV2	793167	985508	69.20%	4.180%
96	14.051	2033	2034	2042	rVB	252415	238865	16.77%	1.013%
97	15.074	2205	2208	2214	rBV	243502	365714	25.68%	1.551%
98	15.380	2257	2260	2266	rVB	168831	205571	14.43%	0.872%
99	15.445	2268	2271	2276	rVB	146977	177421	12.46%	0.753%
100	15.509	2278	2282	2285	rBV	432132	528413	37.10%	2.241%

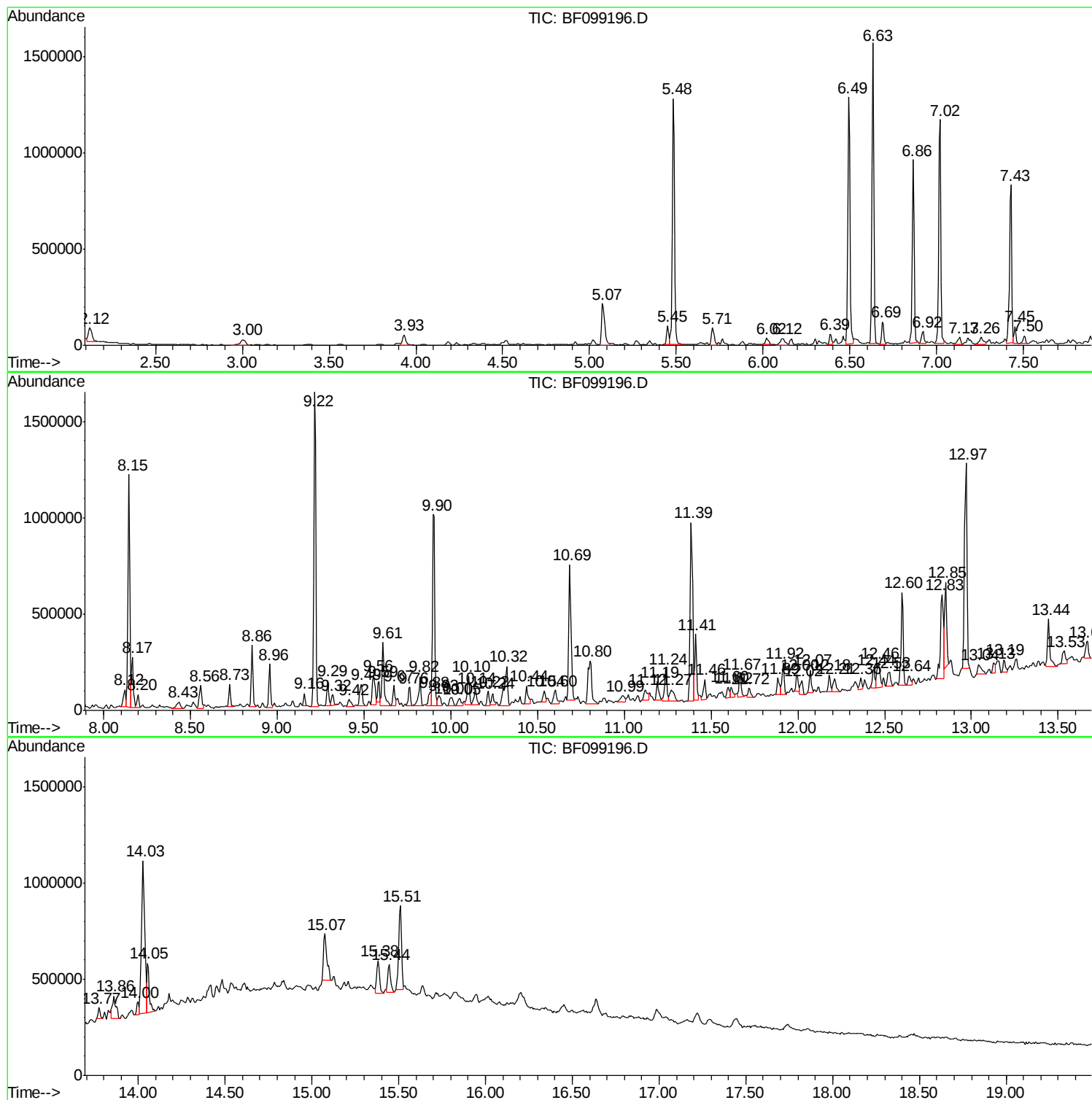
Sum of corrected areas: 23575386

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Instrument :
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ClientSampled :
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF092317.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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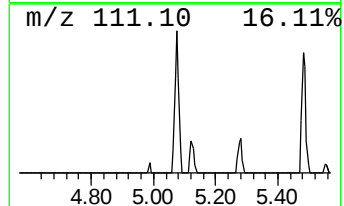
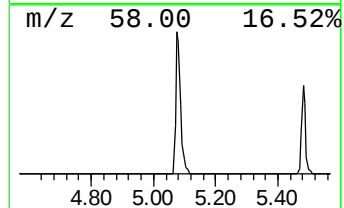
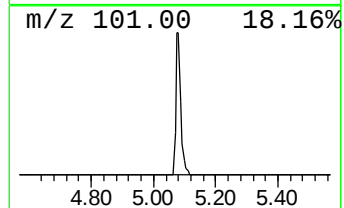
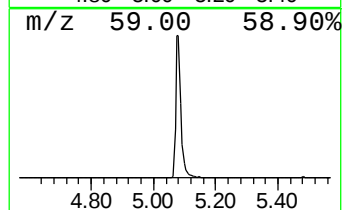
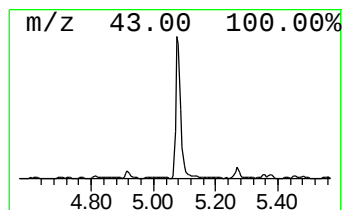
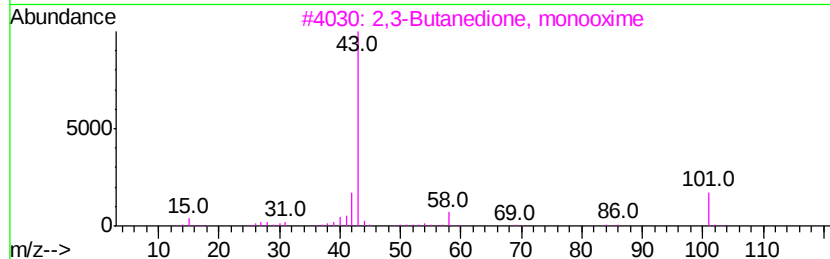
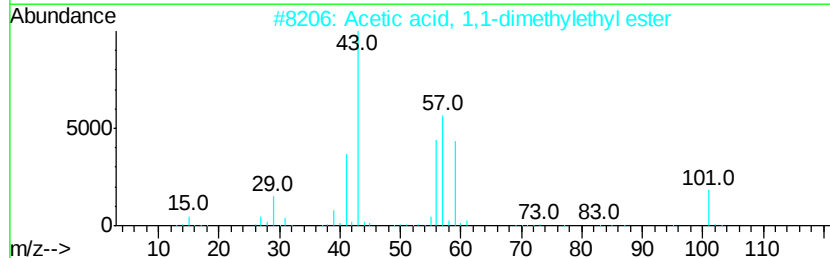
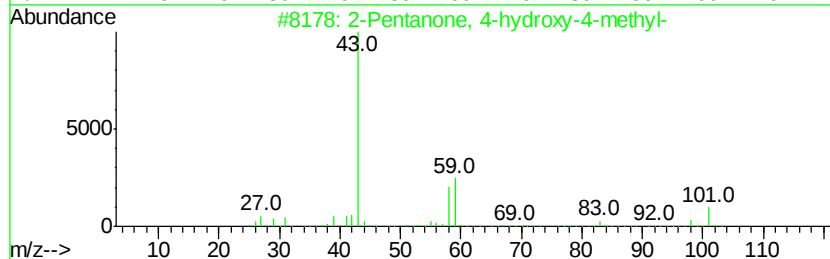
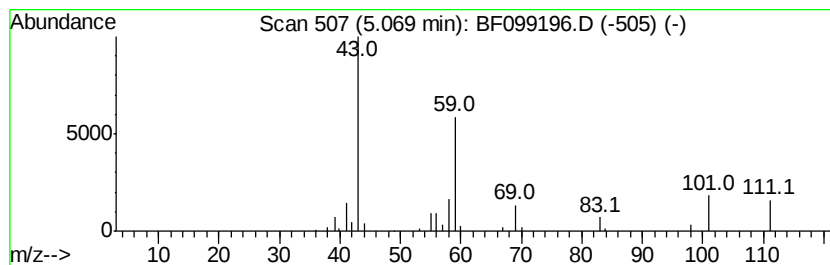
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF092317.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.07	6.16 ng	243427	1,4-Dichlorobenzene-d4	6.86

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	53
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	38
3			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	25
4			3-Hexanol, 4,4-dimethyl-	130	C8H18O	019550-09-5	25
5			n-Propyl t-butyl ether	116	C7H16O	1000334-81-9	25



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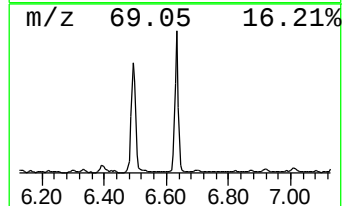
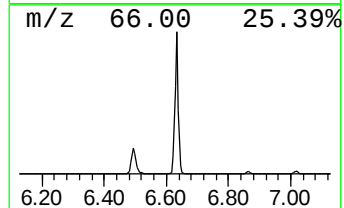
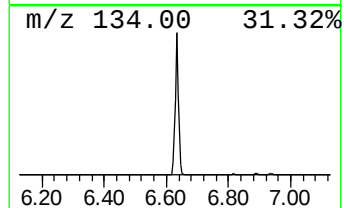
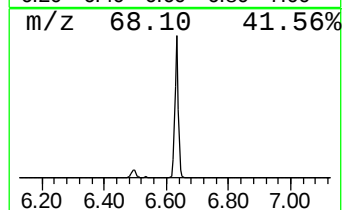
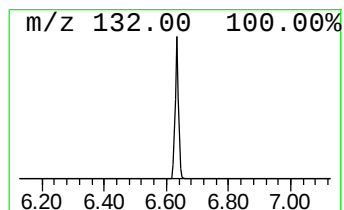
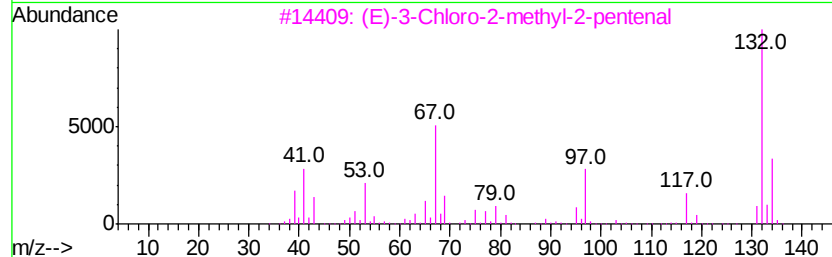
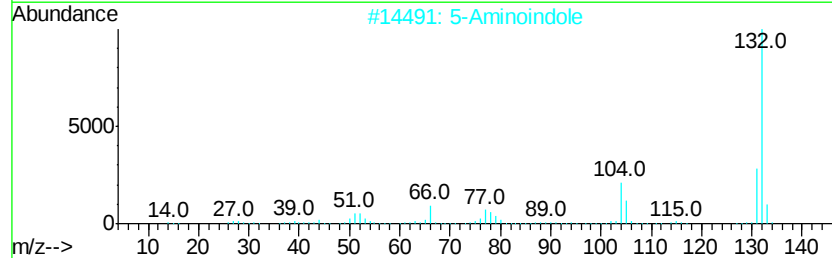
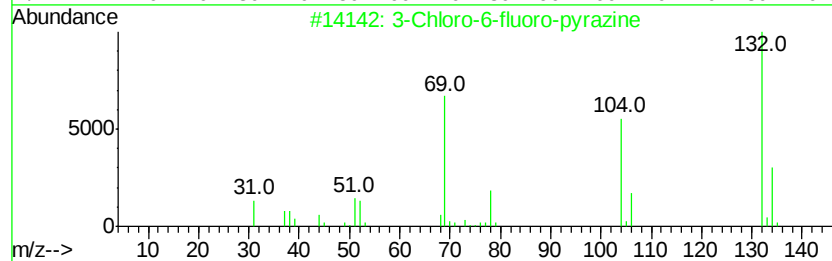
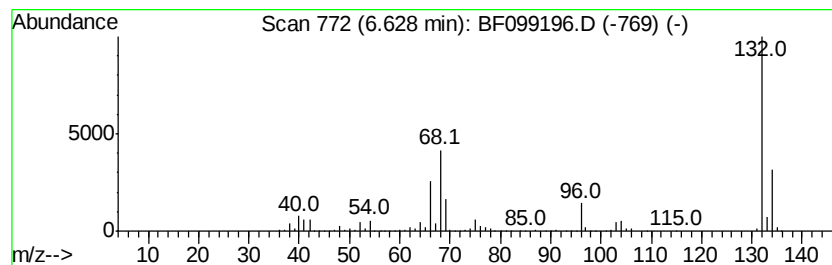
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF092317.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown6.63 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.63	32.27 ng	1274780	1,4-Dichlorobenzene-d4	6.86

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	25
2		5-Aminoindole	132	C8H8N2	005192-03-0	12
3		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	12
4		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9
5		Xenon	132	Xe	007440-63-3	9



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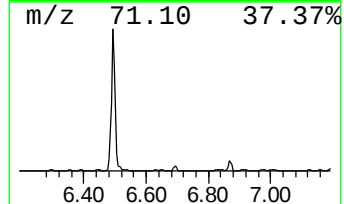
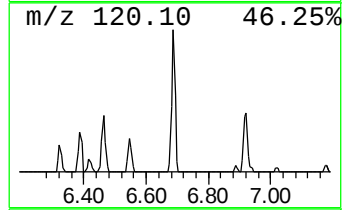
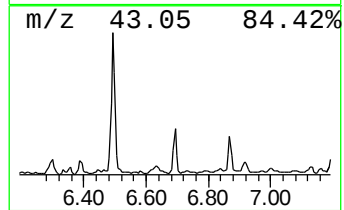
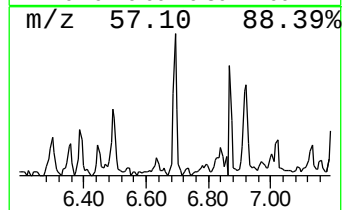
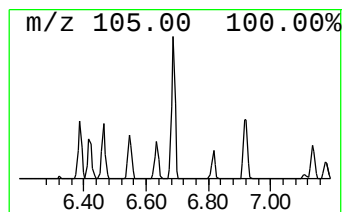
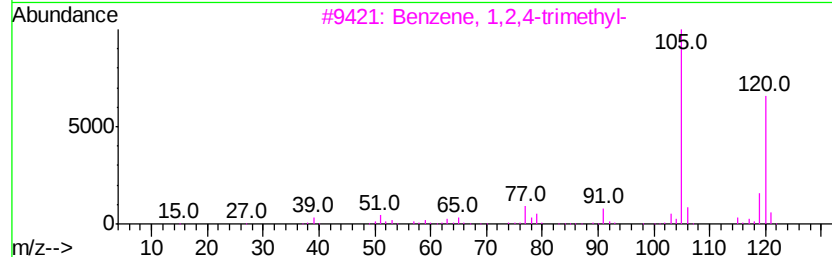
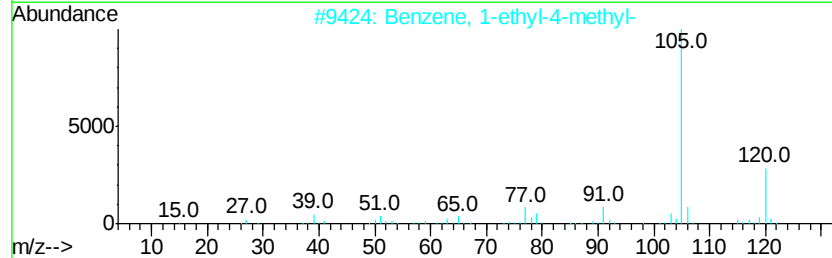
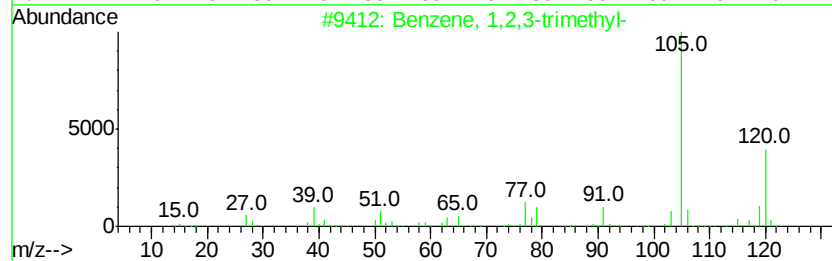
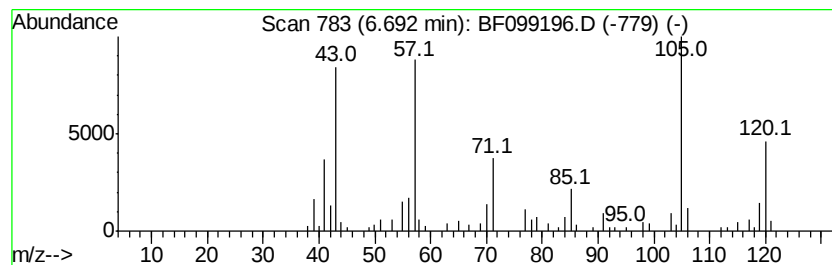
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Benzene, 1,2,3-trimethyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.69	2.88 ng	113886	1,4-Dichlorobenzene-d4	6.86

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	94
2		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	93
3		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
4		Mesitylene	120	C9H12	000108-67-8	91
5		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	90



Data Path : Z:\HPCHEM1\BNA F\DATA\BF100717\
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Acq On : 7 Oct 2017 16:07
Operator : SJ/JU
Sample : I5586-19 2X
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
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ClientSampleId :
G56A-3-6

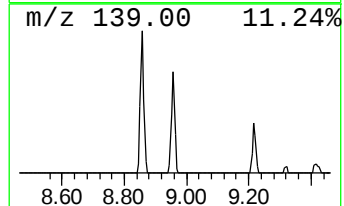
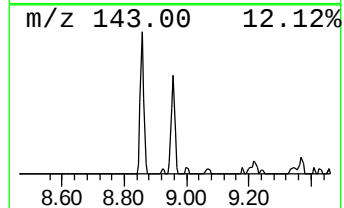
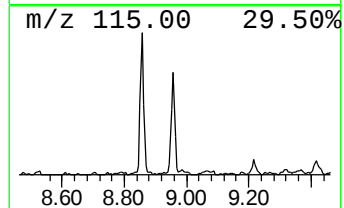
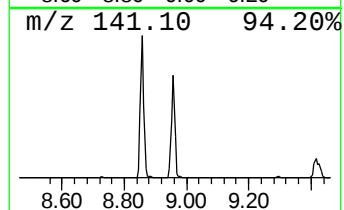
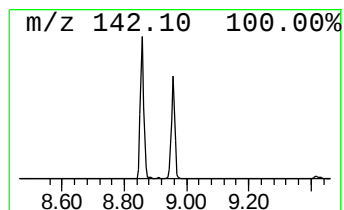
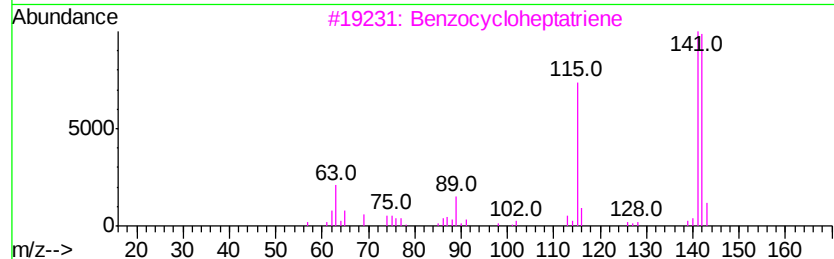
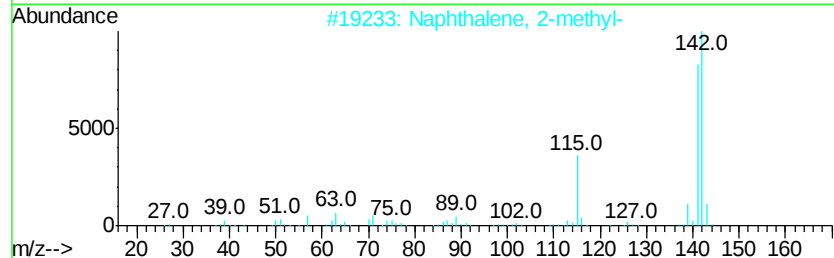
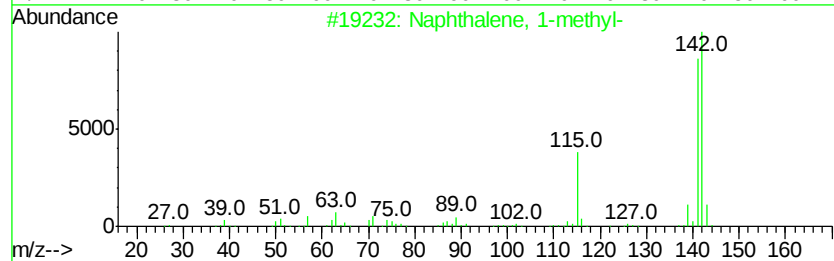
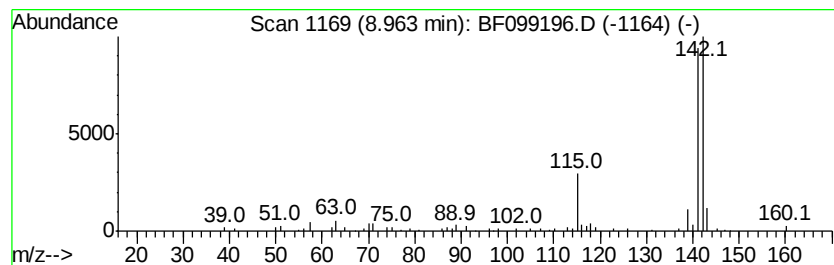
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Naphthalene, 1-methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.96	3.49 ng	171098	Naphthalene-d8	8.15

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
2		Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
3		Benzocycloheptatriene	142	C11H10	000264-09-5	91
4		1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	91
5		1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	64



Data Path : Z:\HPCHEM1\BNA F\DATA\BF100717\
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ClientSampleId :
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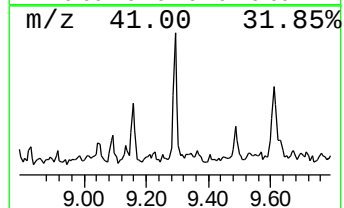
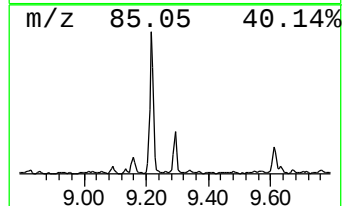
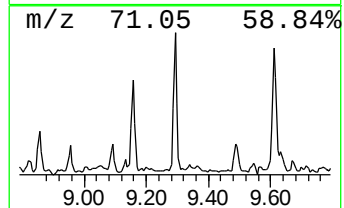
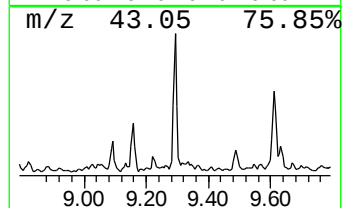
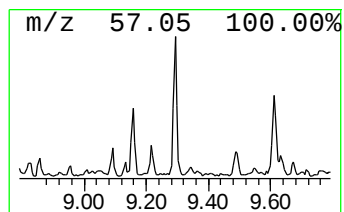
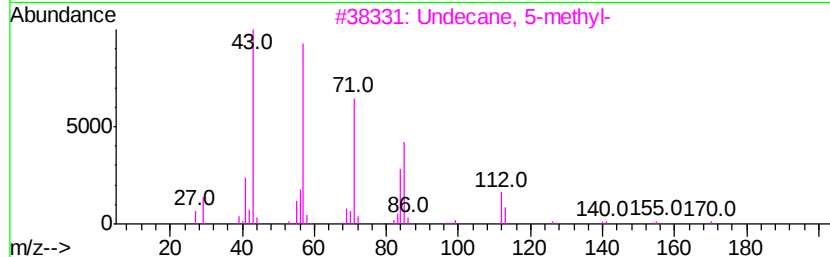
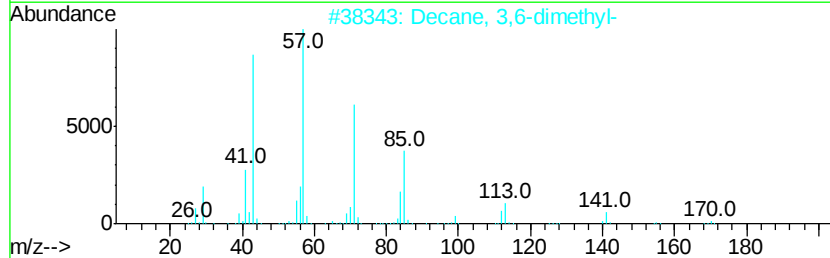
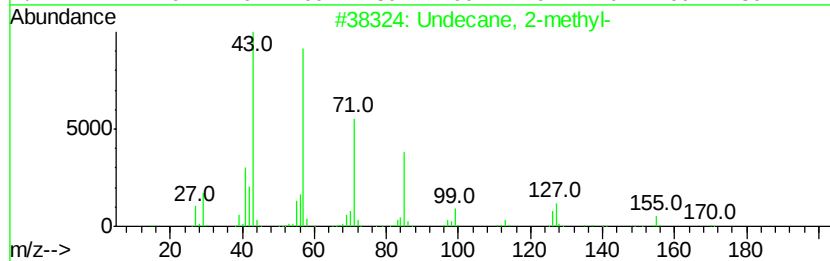
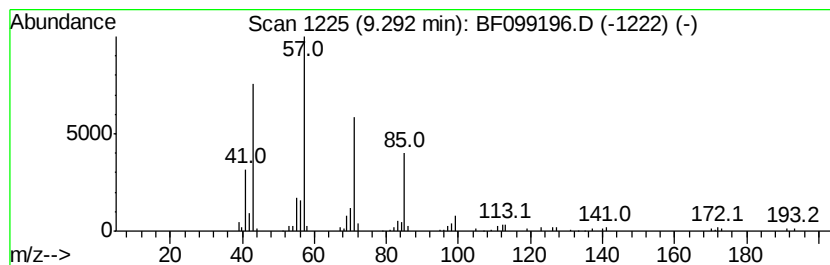
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Undecane, 2-methyl- Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.29	2.42 ng	114026	Acenaphthene-d10	9.90

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Undecane, 2-methyl-	170	C12H26	007045-71-8	87
2		Decane, 3,6-dimethyl-	170	C12H26	017312-53-7	87
3		Undecane, 5-methyl-	170	C12H26	001632-70-8	83
4		Decane, 2,3,6-trimethyl-	184	C13H28	062238-12-4	72
5		Dodecane	170	C12H26	000112-40-3	72



Data Path : Z:\HPCHEM1\BNA F\DATA\BF100717\
Data File : BF099196.D
Acq On : 7 Oct 2017 16:07
Operator : SJ/JU
Sample : I5586-19 2X
Misc :
ALS Vial : 13 Sample Multiplier: 1

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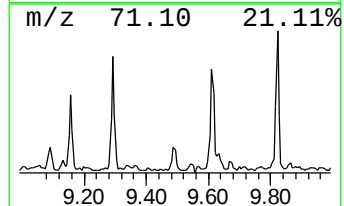
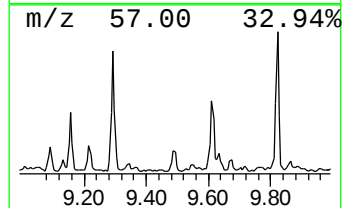
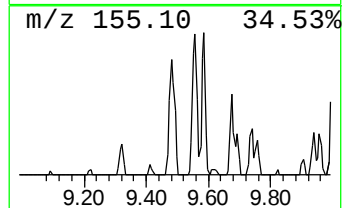
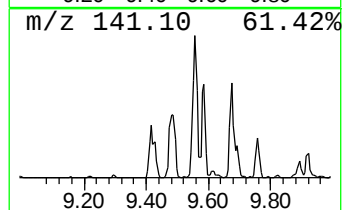
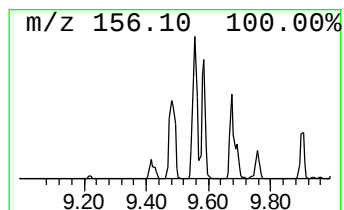
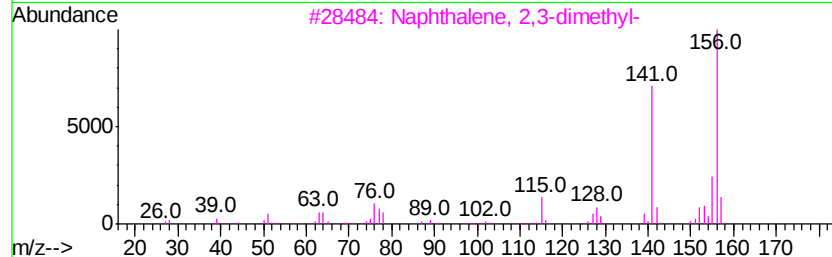
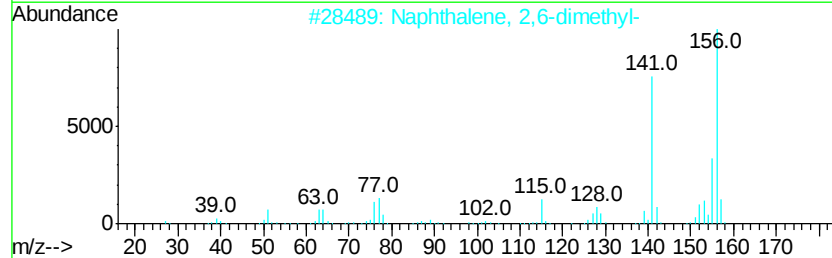
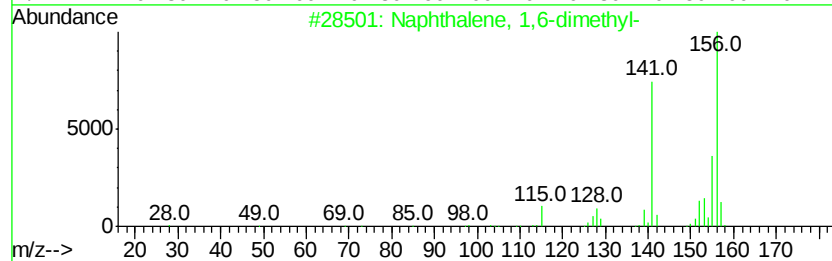
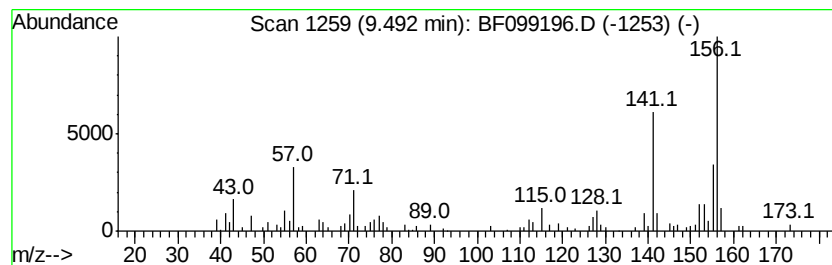
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Naphthalene, 1,6-dimethyl- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.49	2.84 ng	133876	Acenaphthene-d10	9.90

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	96
2		Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	96
3		Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	96
4		Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	96
5		Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	96



Data Path : Z:\HPCHEM1\BNA F\DATA\BF100717\
Data File : BF099196.D
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Misc :
ALS Vial : 13 Sample Multiplier: 1

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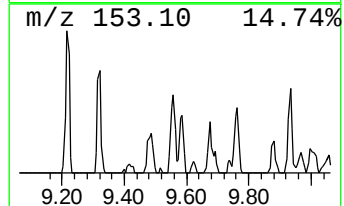
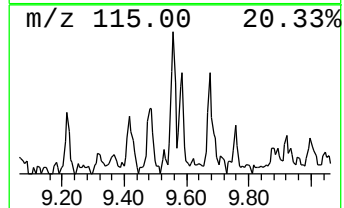
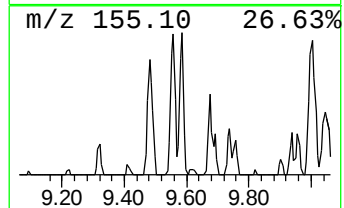
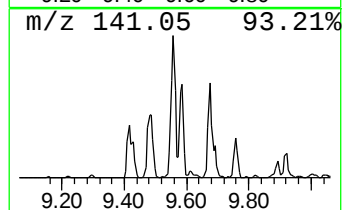
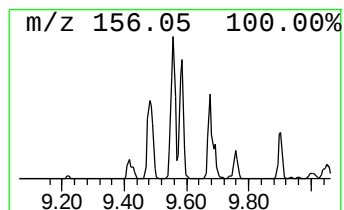
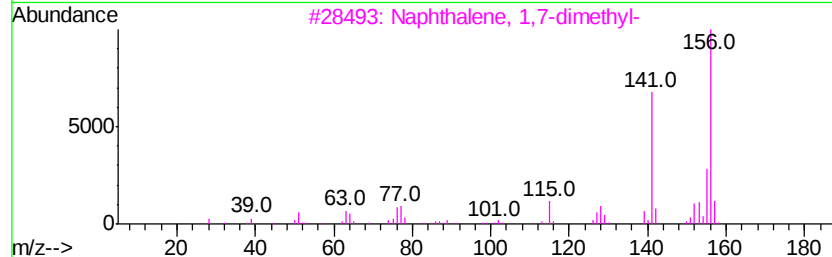
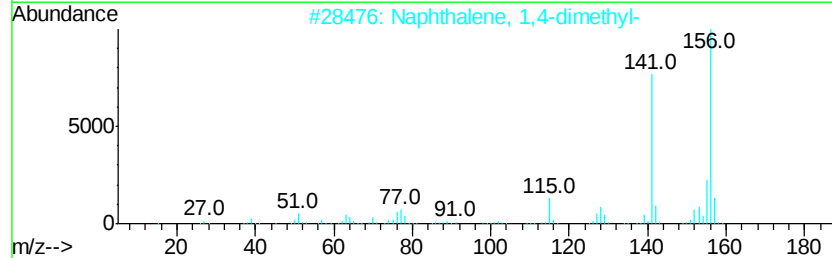
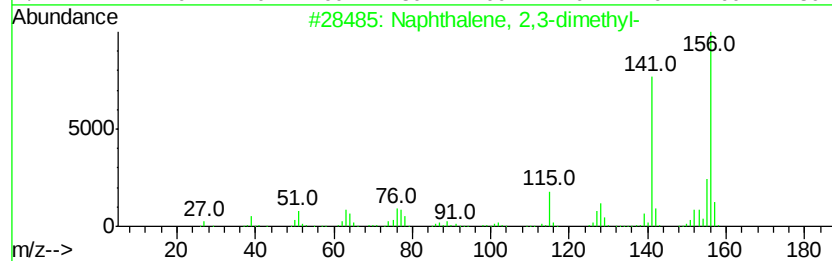
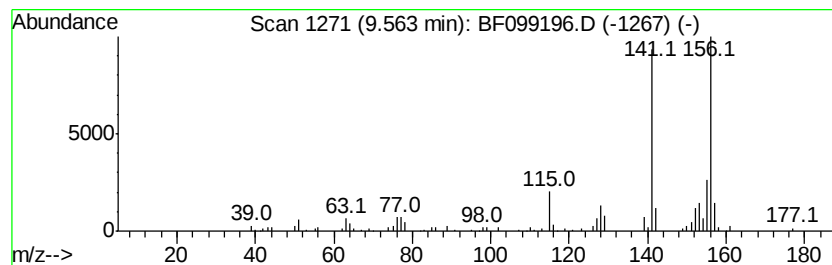
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Naphthalene, 2,3-dimethyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.56	3.19 ng	150399	Acenaphthene-d10	9.90

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
2		Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	97
3		Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	96
4		Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	96
5		Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	96



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF100717\
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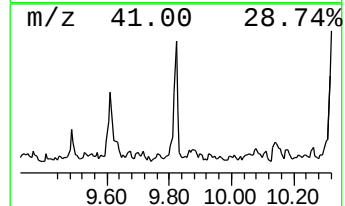
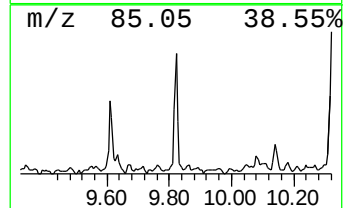
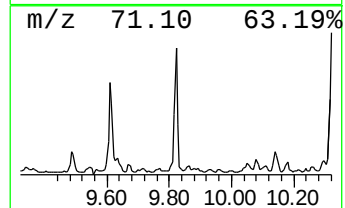
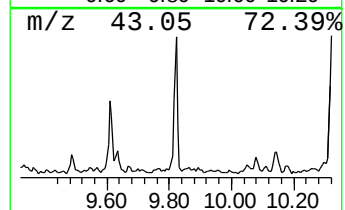
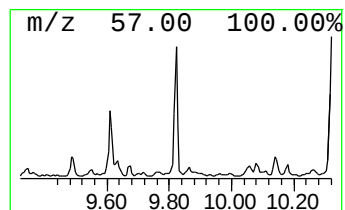
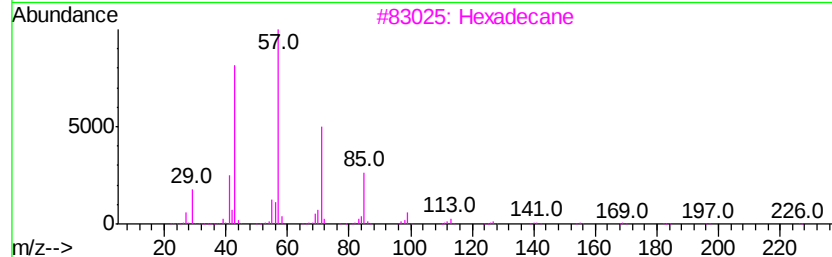
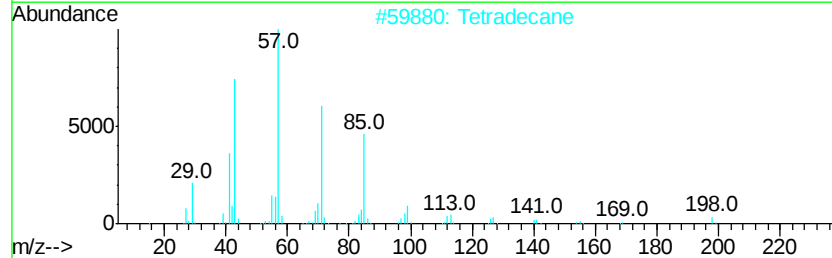
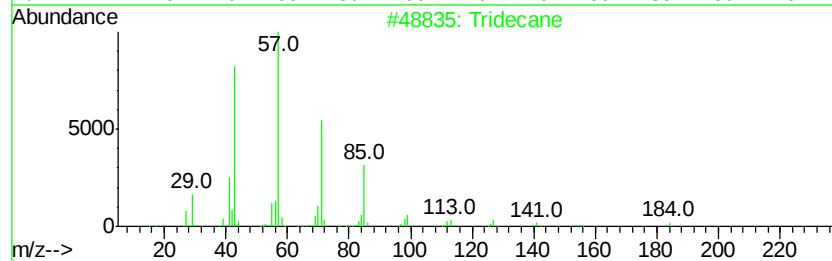
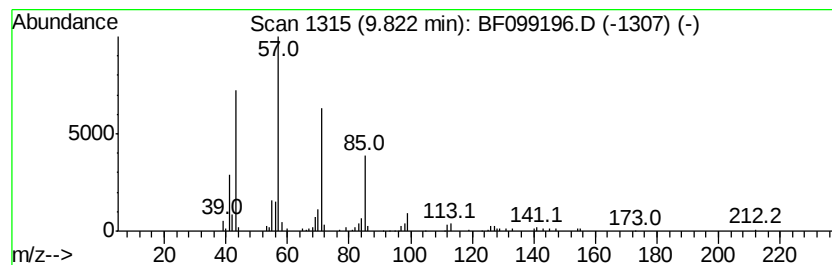
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Tridecane Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.	
9.82	2.70 ng	127241	Acenaphthene-d10	9.90	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tridecane	184	C13H28	000629-50-5	90
2	Tetradecane	198	C14H30	000629-59-4	90
3	Hexadecane	226	C16H34	000544-76-3	80
4	Undecane, 2,10-dimethyl-	184	C13H28	017301-27-8	72
5	Octadecane, 1-chloro-	288	C18H37Cl	003386-33-2	64



Data Path : Z:\HPCHEM1\BNA F\DATA\BF100717\
Data File : BF099196.D
Acq On : 7 Oct 2017 16:07
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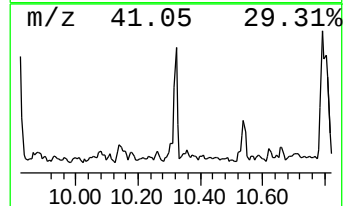
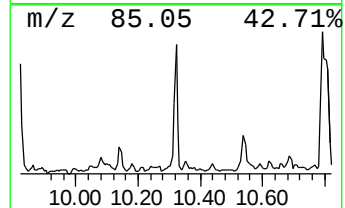
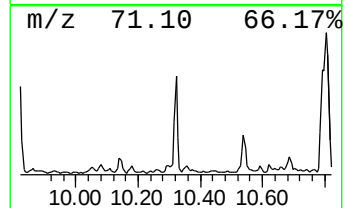
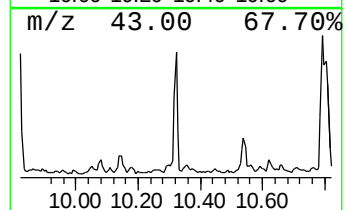
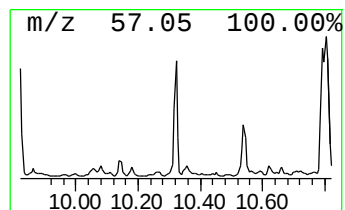
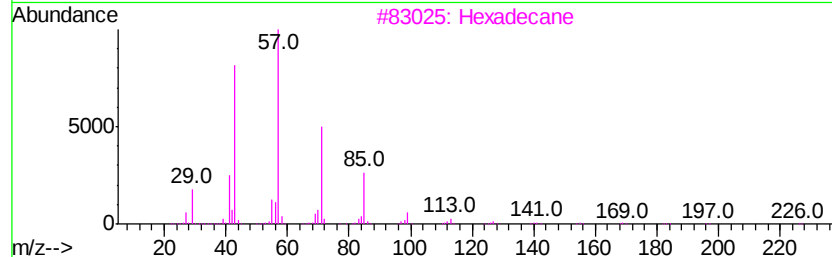
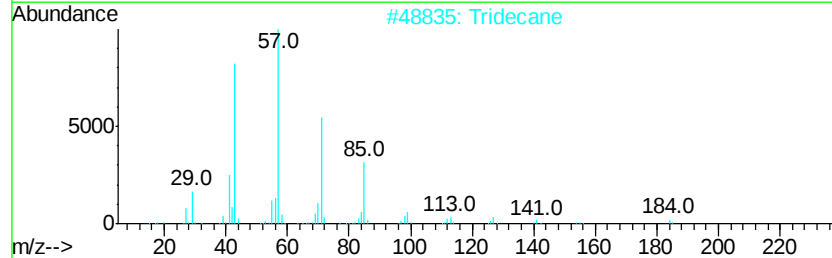
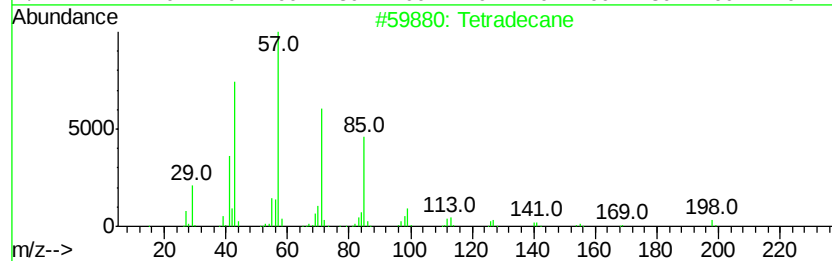
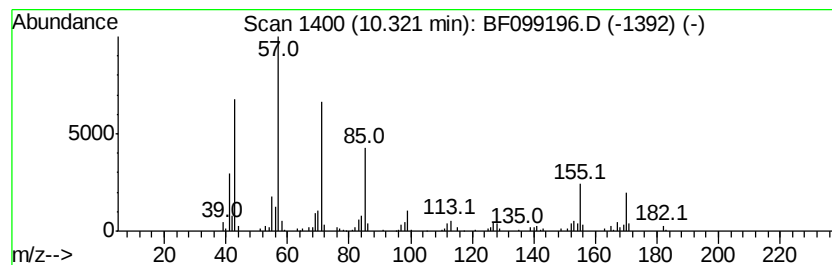
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Tetradecane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.32	5.12 ng	241385	Acenaphthene-d10	9.90

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetradecane	198	C14H30	000629-59-4	58
2			Tridecane	184	C13H28	000629-50-5	52
3			Hexadecane	226	C16H34	000544-76-3	52
4			Sulfurous acid, 2-propyl tetradec...	320	C17H36O3S	1000309-12-5	52
5			Pentadecane	212	C15H32	000629-62-9	52



Data Path : Z:\HPCHEM1\BNA F\DATA\BF100717\
Data File : BF099196.D
Acq On : 7 Oct 2017 16:07
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Misc :
ALS Vial : 13 Sample Multiplier: 1

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ClientSampleId :
G56A-3-6

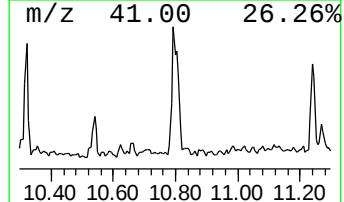
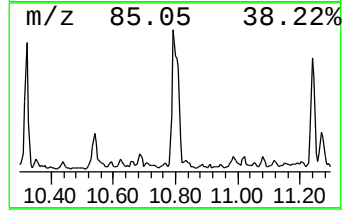
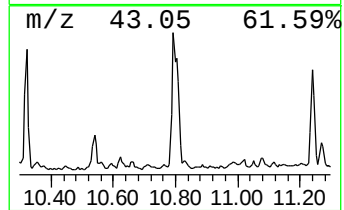
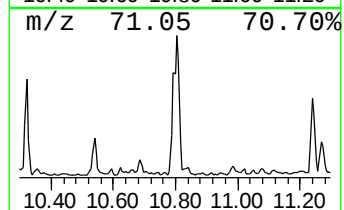
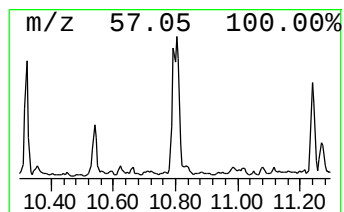
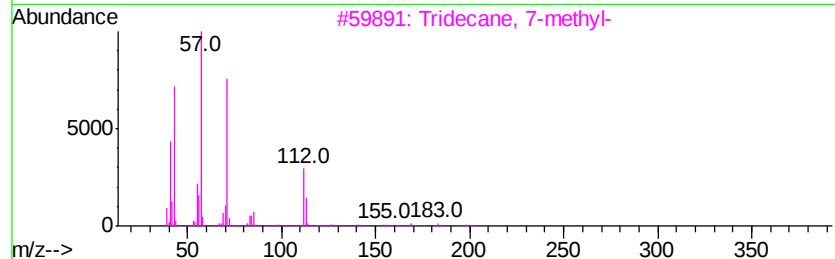
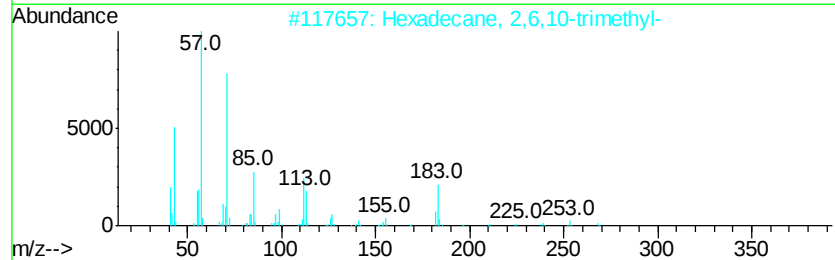
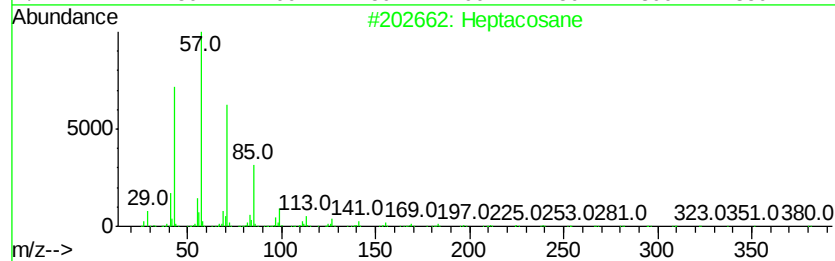
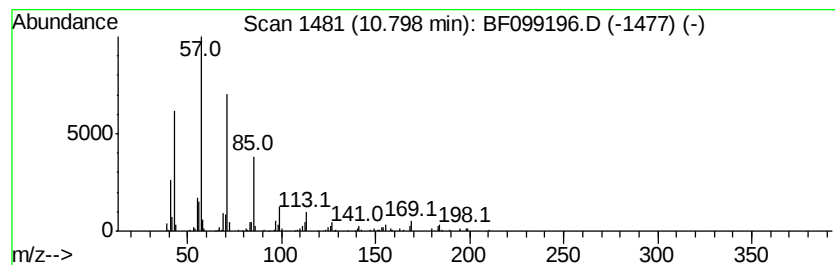
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Heptacosane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.80	9.44 ng	365631	Phenanthrene-d10	11.39

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptacosane	380	C27H56	000593-49-7	80
2			Hexadecane, 2,6,10-trimethyl-	268	C19H40	055000-52-7	80
3			Tridecane, 7-methyl-	198	C14H30	026730-14-3	76
4			Dodecane, 2,7,10-trimethyl-	212	C15H32	074645-98-0	72
5			Pentadecane, 2,6,10,14-tetramethyl-	268	C19H40	001921-70-6	72



Data Path : Z:\HPCHEM1\BNA F\DATA\BF100717\
Data File : BF099196.D
Acq On : 7 Oct 2017 16:07
Operator : SJ/JU
Sample : I5586-19 2X
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
G56A-3-6

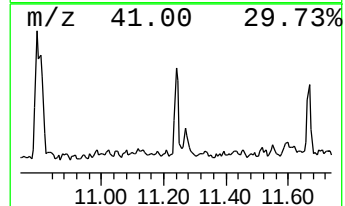
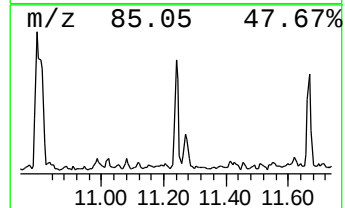
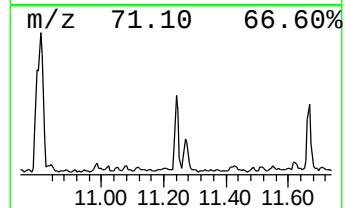
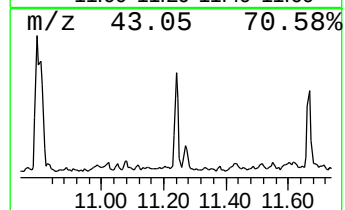
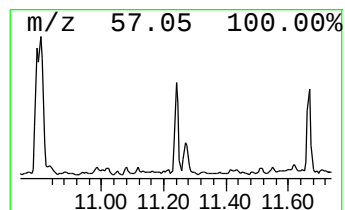
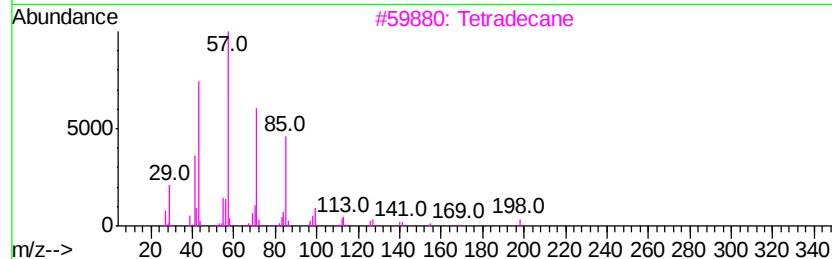
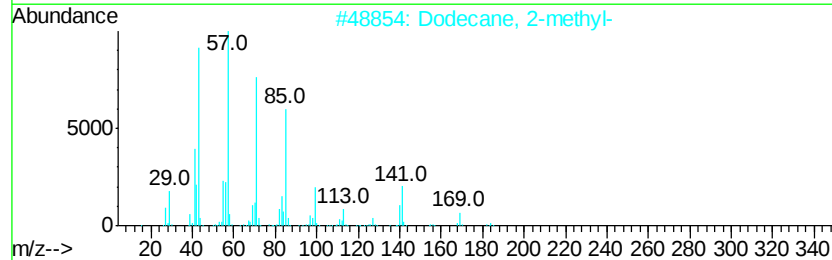
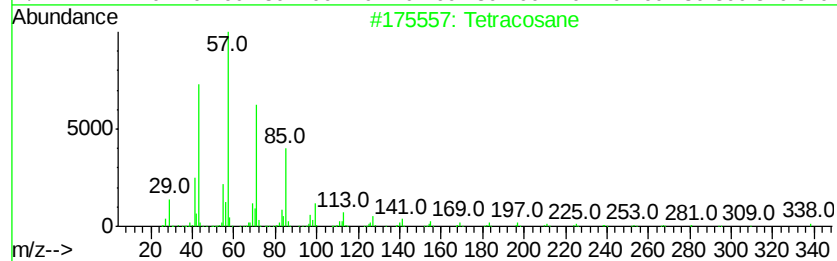
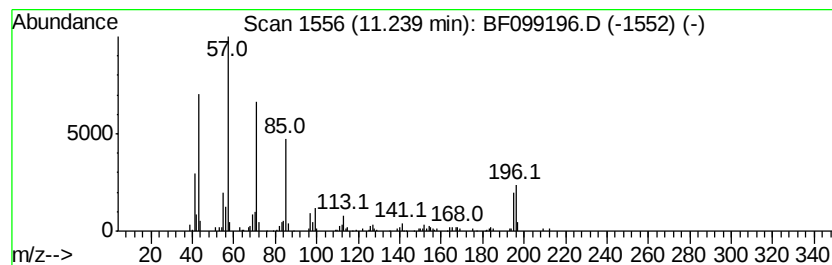
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Tetracosane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.24	4.34 ng	168302	Phenanthrene-d10	11.39

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetracosane	338	C24H50	000646-31-1	58
2		Dodecane, 2-methyl-	184	C13H28	001560-97-0	53
3		Tetradecane	198	C14H30	000629-59-4	52
4		Dodecane, 1-iodo-	296	C12H25I	004292-19-7	52
5		Decane, 3-bromo-	220	C10H21Br	030571-71-2	49



Data Path : Z:\HPCHEM1\BNA F\DATA\BF100717\
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ClientSampleId :
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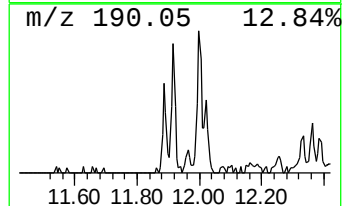
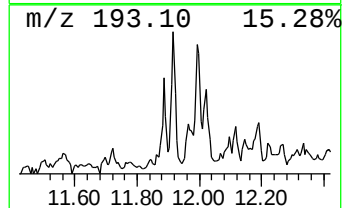
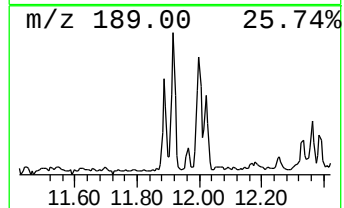
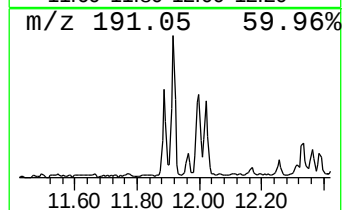
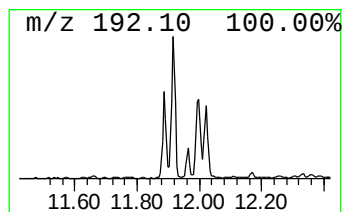
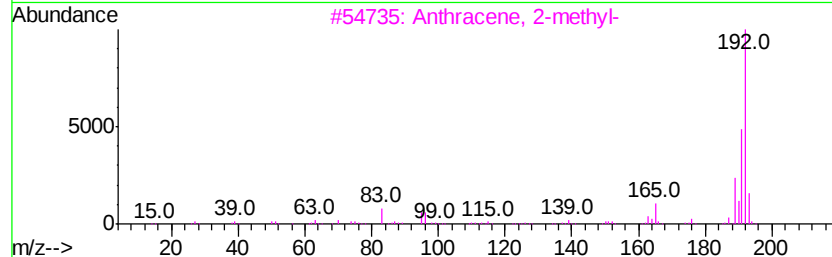
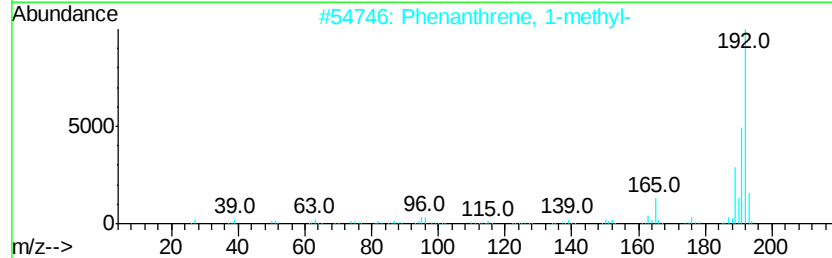
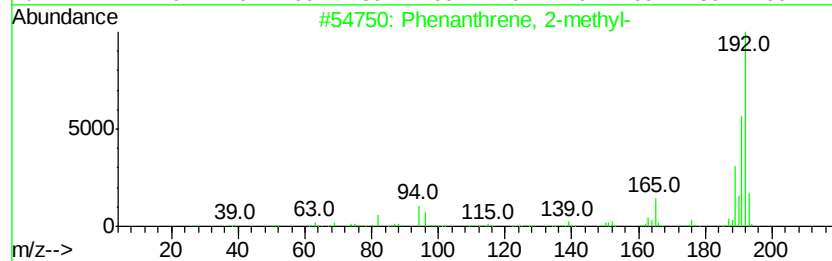
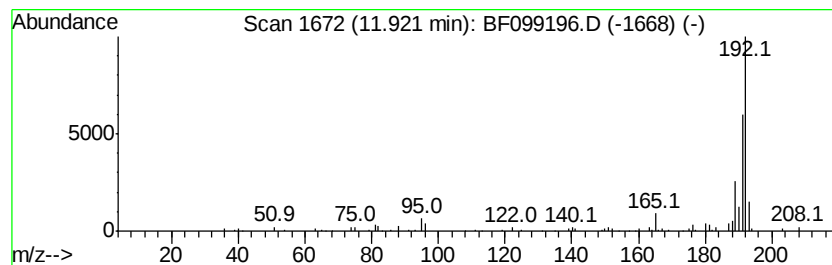
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Phenanthrene, 2-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.92	3.90 ng	151091	Phenanthrene-d10	11.39

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	98
2		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	97
3		Anthracene, 2-methyl-	192	C15H12	000613-12-7	96
4		Anthracene, 9-methyl-	192	C15H12	000779-02-2	96
5		1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	95



Data Path : Z:\HPCHEM1\BNA F\DATA\BF100717\
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Acq On : 7 Oct 2017 16:07
Operator : SJ/JU
Sample : I5586-19 2X
Misc :
ALS Vial : 13 Sample Multiplier: 1

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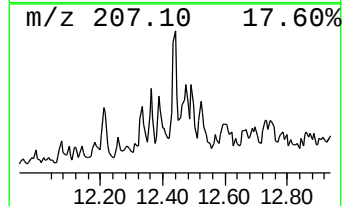
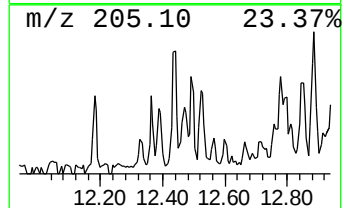
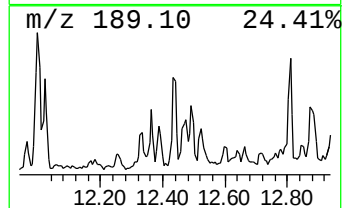
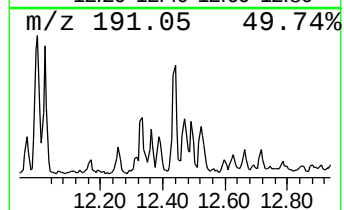
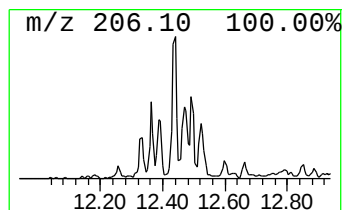
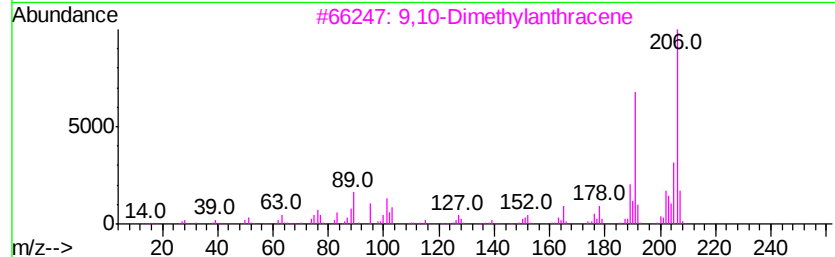
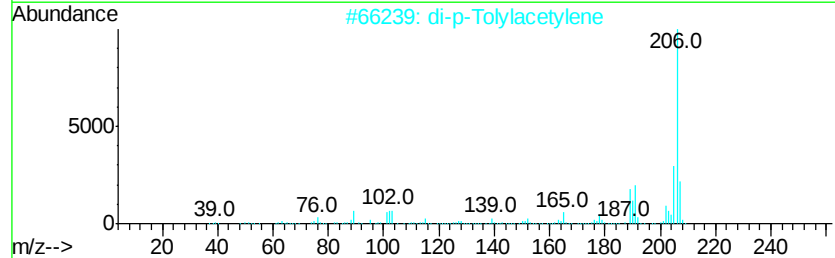
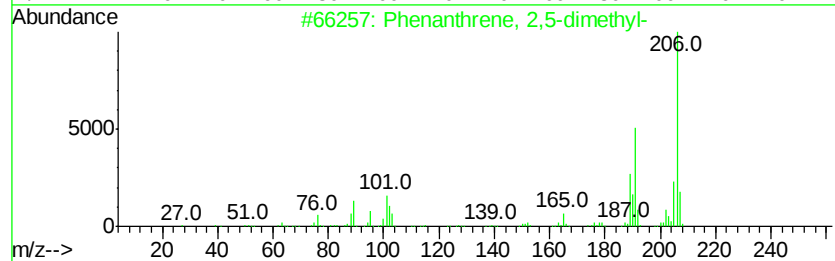
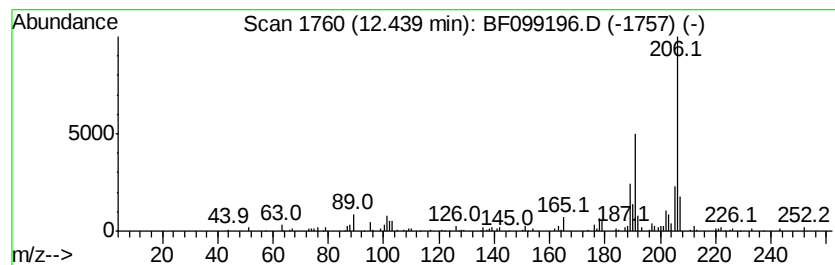
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 Phenanthrene, 2,5-dimethyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.44	2.87 ng	111076	Phenanthrene-d10	11.39

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	96
2		di-p-Tolylacetylene	206	C16H14	002789-88-0	92
3		9,10-Dimethylanthracene	206	C16H14	000781-43-1	91
4		1H-Indene, 2-methyl-3-phenyl-	206	C16H14	035099-60-6	90
5		Phenanthrene, 1,7-dimethyl-	206	C16H14	000483-87-4	90



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF100717\
Data File : BF099196.D
Acq On : 7 Oct 2017 16:07
Operator : SJ/JU
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Misc :
ALS Vial : 13 Sample Multiplier: 1

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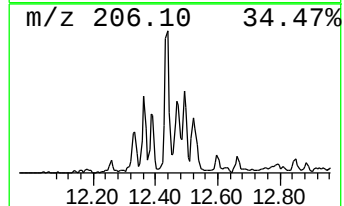
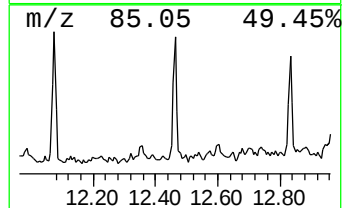
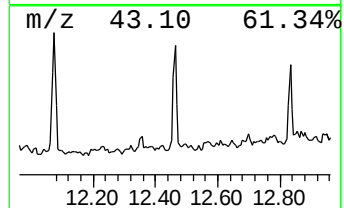
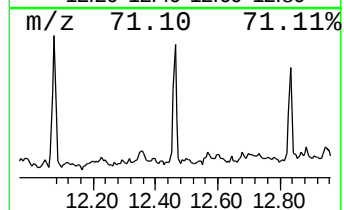
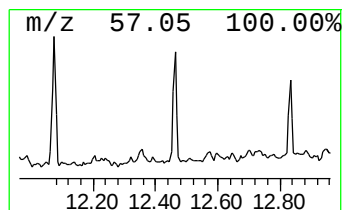
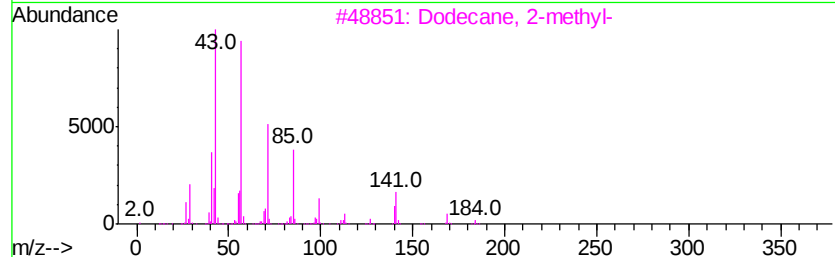
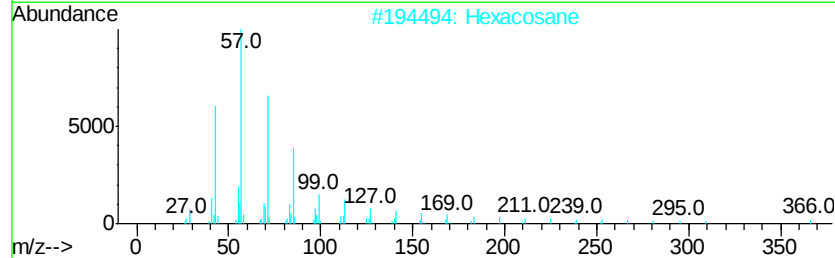
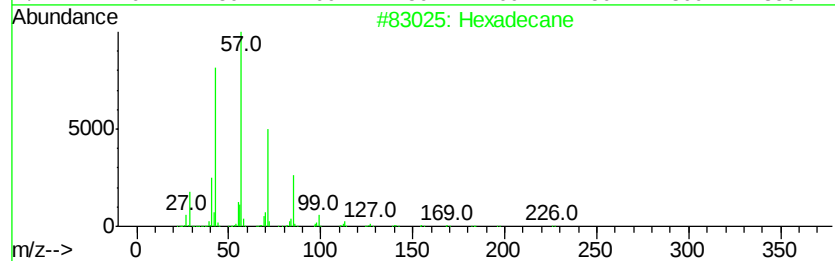
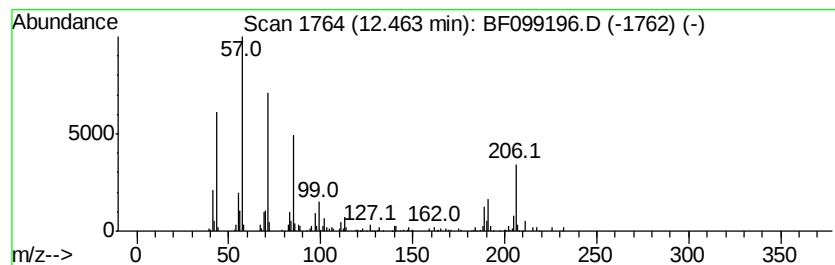
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 Hexadecane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.46	3.52 ng	136356	Phenanthrene-d10	11.39

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane	226	C16H34	000544-76-3	58
2	Hexacosane	366	C26H54	000630-01-3	53
3	Dodecane, 2-methyl-	184	C13H28	001560-97-0	52
4	Octacosane	394	C28H58	000630-02-4	49
5	Nonane, 3,7-dimethyl-	156	C11H24	017302-32-8	47



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF100717\
Data File : BF099196.D
Acq On : 7 Oct 2017 16:07
Operator : SJ/JU
Sample : I5586-19 2X
Misc :
ALS Vial : 13 Sample Multiplier: 1

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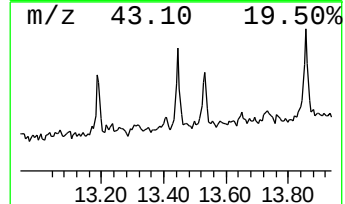
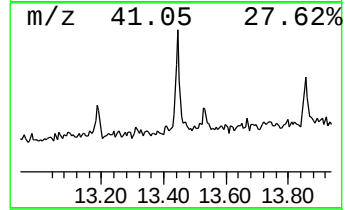
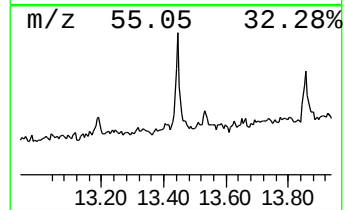
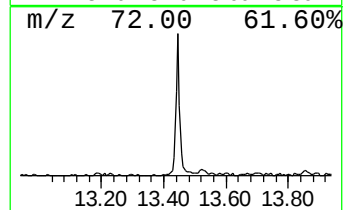
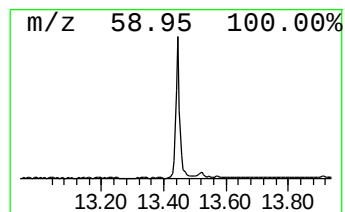
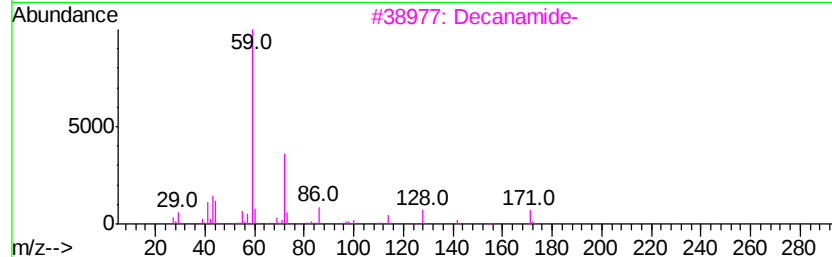
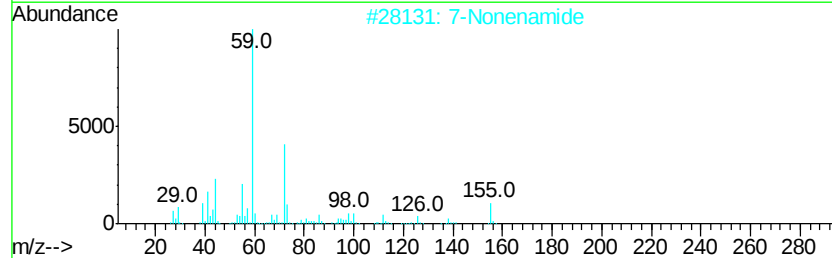
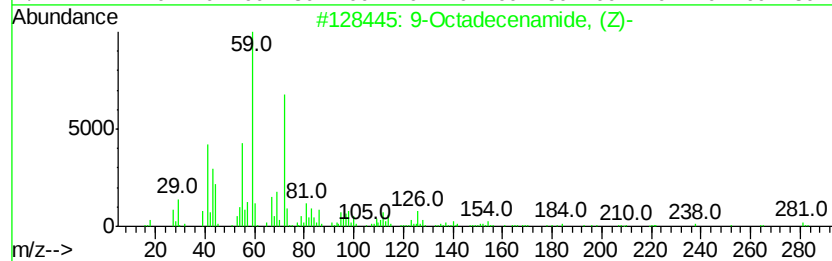
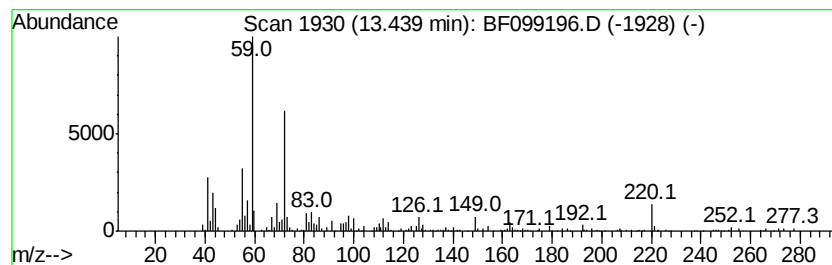
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 9-Octadecenamide, (Z)- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.44	4.81 ng	236778	Chrysene-d12	14.03

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	93
2		7-Nonenamide	155	C9H17NO	090949-53-4	72
3		Decanamide-	171	C10H21NO	002319-29-1	72
4		Pentadecanamide, 15-bromo-	319	C15H30BrNO	1000163-86-1	53
5		Nonanamide	157	C9H19NO	001120-07-6	50



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF100717\
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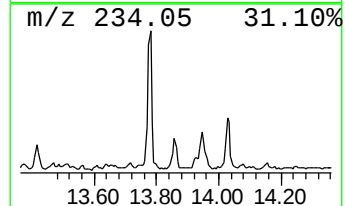
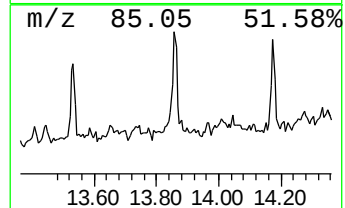
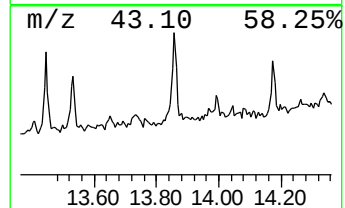
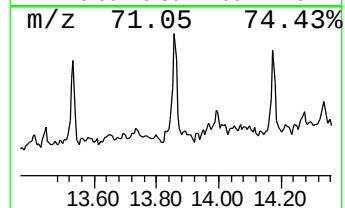
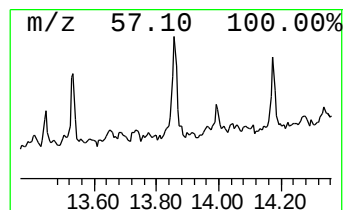
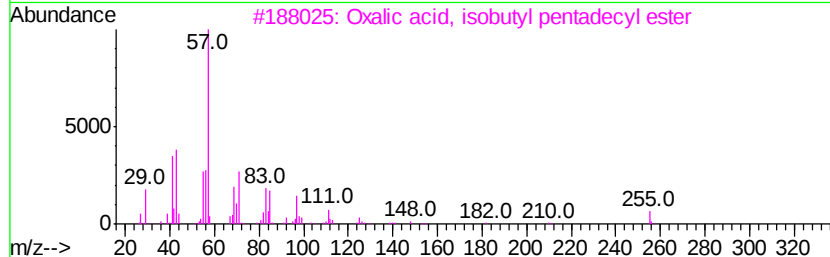
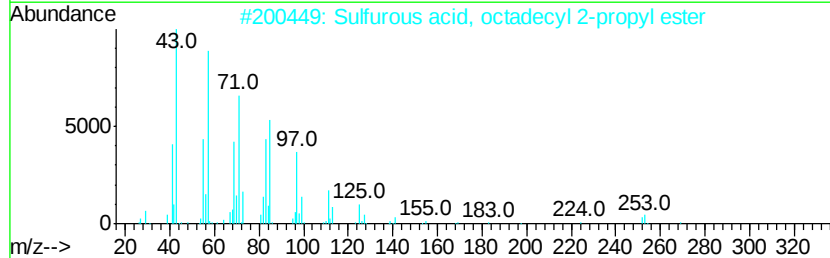
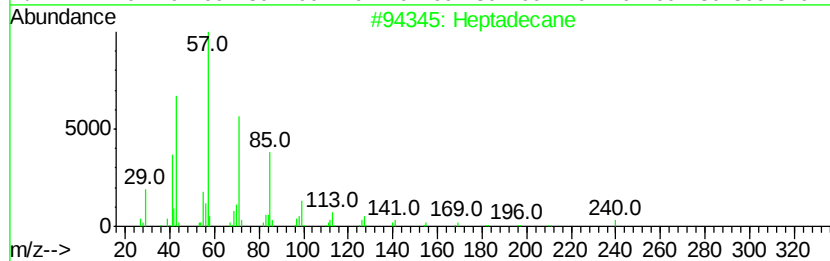
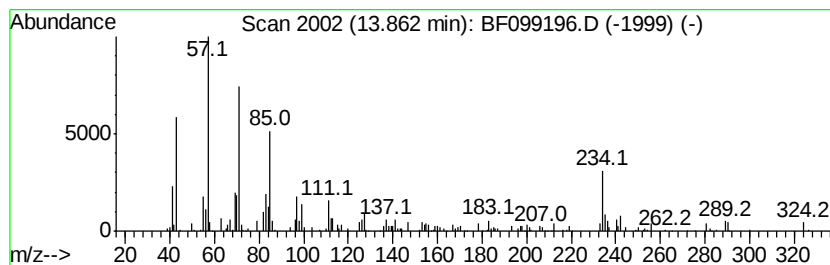
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 19 Heptadecane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.86	3.12 ng	153573	Chrysene-d12	14.03	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane	240	C17H36	000629-78-7	90
2	Sulfurous acid, octadecyl 2-prop...	376	C21H44O3S	1000309-12-7	68
3	Oxalic acid, isobutyl pentadecyl...	356	C21H40O4	1000309-38-0	64
4	Ethanol, 2-(octadecyloxy)-	314	C20H42O2	002136-72-3	62
5	Tridecane	184	C13H28	000629-50-5	55



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF100717\
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 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
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 ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF092317.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
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unknown6.63	6.63	32.3	ng	1274780	1	6.86	790041	20.0
Benzene, 1,2,3-tr...	6.69	2.9	ng	113886	1	6.86	790041	20.0
Naphthalene, 1-me...	8.96	3.5	ng	171098	2	8.15	981512	20.0
Undecane, 2-methyl-	9.29	2.4	ng	114026	3	9.90	942542	20.0
Naphthalene, 1,6-...	9.49	2.8	ng	133876	3	9.90	942542	20.0
Naphthalene, 2,3-...	9.56	3.2	ng	150399	3	9.90	942542	20.0
Tridecane	9.82	2.7	ng	127241	3	9.90	942542	20.0
Tetradecane	10.32	5.1	ng	241385	3	9.90	942542	20.0
Heptacosane	10.80	9.4	ng	365631	4	11.39	774742	20.0
Tetracosane	11.24	4.3	ng	168302	4	11.39	774742	20.0
Phenanthrene, 2-m...	11.92	3.9	ng	151091	4	11.39	774742	20.0
Phenanthrene, 2,5...	12.44	2.9	ng	111076	4	11.39	774742	20.0
Hexadecane	12.46	3.5	ng	136356	4	11.39	774742	20.0
9-Octadecenamide, ...	13.44	4.8	ng	236778	5	14.03	985508	20.0
Heptadecane	13.86	3.1	ng	153573	5	14.03	985508	20.0