

Data Path : Z:\HPCHEM1\BNA F\DATA\BF111717\
 Data File : BF100669.D
 Acq On : 17 Nov 2017 11:41
 Operator : SJ/JU
 Sample : I6477-01 10X
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TOTE-1-2

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-BF110817.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	2.287	31	34	42	rVB3	8702	15501	1.44%	0.113%
2	3.963	313	319	325	rBV	171650	220310	20.42%	1.605%
3	5.487	575	578	585	rBV	126974	105488	9.78%	0.769%
4	6.492	737	749	753	rBV2	89175	124323	11.53%	0.906%
5	6.645	771	775	778	rBV	272656	233722	21.67%	1.703%
6	6.881	811	815	818	rBV	744855	598821	55.51%	4.363%
7	6.933	821	824	827	rVV	113682	93297	8.65%	0.680%
8	6.969	827	830	835	rVV2	16153	18322	1.70%	0.133%
9	7.033	835	841	845	rVV	288308	251425	23.31%	1.832%
10	7.222	865	873	876	rBV	37072	53005	4.91%	0.386%
11	7.433	906	909	913	rBV	190188	166602	15.44%	1.214%
12	7.528	920	925	927	rBV2	21790	23462	2.17%	0.171%
13	7.569	927	932	938	rVB4	26437	38618	3.58%	0.281%
14	7.657	944	947	950	rBV	47993	50335	4.67%	0.367%
15	7.686	950	952	954	rVV	47126	41650	3.86%	0.303%
16	7.704	954	955	958	rVB2	18821	13679	1.27%	0.100%
17	7.822	971	975	980	rBV3	19083	28398	2.63%	0.207%
18	7.892	980	987	990	rVV2	53300	63049	5.84%	0.459%
19	7.951	993	997	1005	rVV	193625	176629	16.37%	1.287%
20	8.063	1012	1016	1019	rBV3	21626	23729	2.20%	0.173%
21	8.133	1024	1028	1029	rBV3	11298	14582	1.35%	0.106%
22	8.163	1029	1033	1036	rVV	1003211	818379	75.87%	5.963%
23	8.233	1042	1045	1049	rVB2	27991	24673	2.29%	0.180%
24	8.351	1060	1065	1068	rBV5	18153	27170	2.52%	0.198%
25	8.439	1076	1080	1083	rBV3	21893	22573	2.09%	0.164%
26	8.475	1083	1086	1089	rBV	36724	45110	4.18%	0.329%
27	8.510	1089	1092	1094	rBV	108212	97302	9.02%	0.709%
28	8.586	1102	1105	1109	rVB5	17022	22655	2.10%	0.165%
29	8.686	1117	1122	1123	rBV3	9955	15944	1.48%	0.116%
30	8.727	1124	1129	1131	rVV3	14027	19127	1.77%	0.139%
31	8.769	1131	1136	1141	rVB9	11246	20954	1.94%	0.153%
32	8.839	1141	1148	1153	rBV6	34778	69815	6.47%	0.509%
33	8.886	1153	1156	1160	rVV5	11415	15508	1.44%	0.113%
34	8.957	1166	1168	1172	rVB4	20751	17932	1.66%	0.131%

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35	9.039	1180	1182	1185	rBV3	49224	42164	3.91%	0.307%
36	9.075	1185	1188	1195	rVB3	63851	75620	7.01%	0.551%
37	9.175	1204	1205	1209	rVB3	20467	18398	1.71%	0.134%
38	9.233	1211	1215	1219	rVB	517832	403682	37.42%	2.941%
39	9.310	1224	1228	1229	rBV4	15703	19250	1.78%	0.140%
40	9.327	1229	1231	1236	rVB4	27574	29617	2.75%	0.216%
41	9.386	1236	1241	1244	rBV	296163	262206	24.31%	1.910%
42	9.469	1253	1255	1257	rVB2	19103	14539	1.35%	0.106%
43	9.498	1257	1260	1263	rBV5	20304	28128	2.61%	0.205%
44	9.533	1263	1266	1269	rBV3	17302	20158	1.87%	0.147%
45	9.569	1269	1272	1274	rBV3	24905	28378	2.63%	0.207%
46	9.669	1285	1289	1292	rBV6	24595	39058	3.62%	0.285%
47	9.822	1310	1315	1317	rBV4	38347	61907	5.74%	0.451%
48	9.880	1323	1325	1327	rBV	25112	23461	2.17%	0.171%
49	9.916	1327	1331	1335	rVV	976243	920027	85.29%	6.703%
50	10.127	1364	1367	1371	rBV	146830	131838	12.22%	0.961%
51	10.322	1398	1400	1401	rBV2	24865	19016	1.76%	0.139%
52	10.363	1405	1407	1412	rVV4	32135	31502	2.92%	0.230%
53	10.457	1420	1423	1427	rBV2	89500	79238	7.35%	0.577%
54	10.704	1461	1465	1471	rVB	237723	237074	21.98%	1.727%
55	10.774	1474	1477	1481	rBV	771989	635560	58.92%	4.631%
56	10.816	1481	1484	1488	rVB5	46474	73833	6.84%	0.538%
57	11.169	1542	1544	1547	rBV	53643	41551	3.85%	0.303%
58	11.404	1580	1584	1591	rBV	1086929	931472	86.35%	6.787%
59	11.598	1614	1617	1622	rBV4	37125	55824	5.18%	0.407%
60	11.680	1628	1631	1635	rBV	133010	115786	10.73%	0.844%
61	11.968	1676	1680	1689	rVB	1207045	1038317	96.26%	7.565%
62	12.751	1811	1813	1820	rVB	114813	105084	9.74%	0.766%
63	12.980	1849	1852	1855	rBV	380808	333682	30.93%	2.431%
64	13.015	1855	1858	1872	rVB	1187709	1078713	100.00%	7.860%
65	13.704	1972	1975	1981	rVB	72563	75912	7.04%	0.553%
66	13.939	2012	2015	2024	rVB	835244	823694	76.36%	6.001%
67	14.009	2024	2027	2029	rVV	162791	140990	13.07%	1.027%
68	14.039	2029	2032	2036	rVB	677183	612916	56.82%	4.466%
69	14.557	2117	2120	2124	rBV	39347	49141	4.56%	0.358%
70	14.786	2155	2159	2175	rVB	521543	633220	58.70%	4.614%
71	15.515	2278	2283	2288	rBV	534905	660979	61.27%	4.816%

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Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
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72	15.809	2328	2333	2341	rBV2	127824	228811	21.21%	1.667%
73	16.474	2443	2446	2454	rVB5	17392	32086	2.97%	0.234%

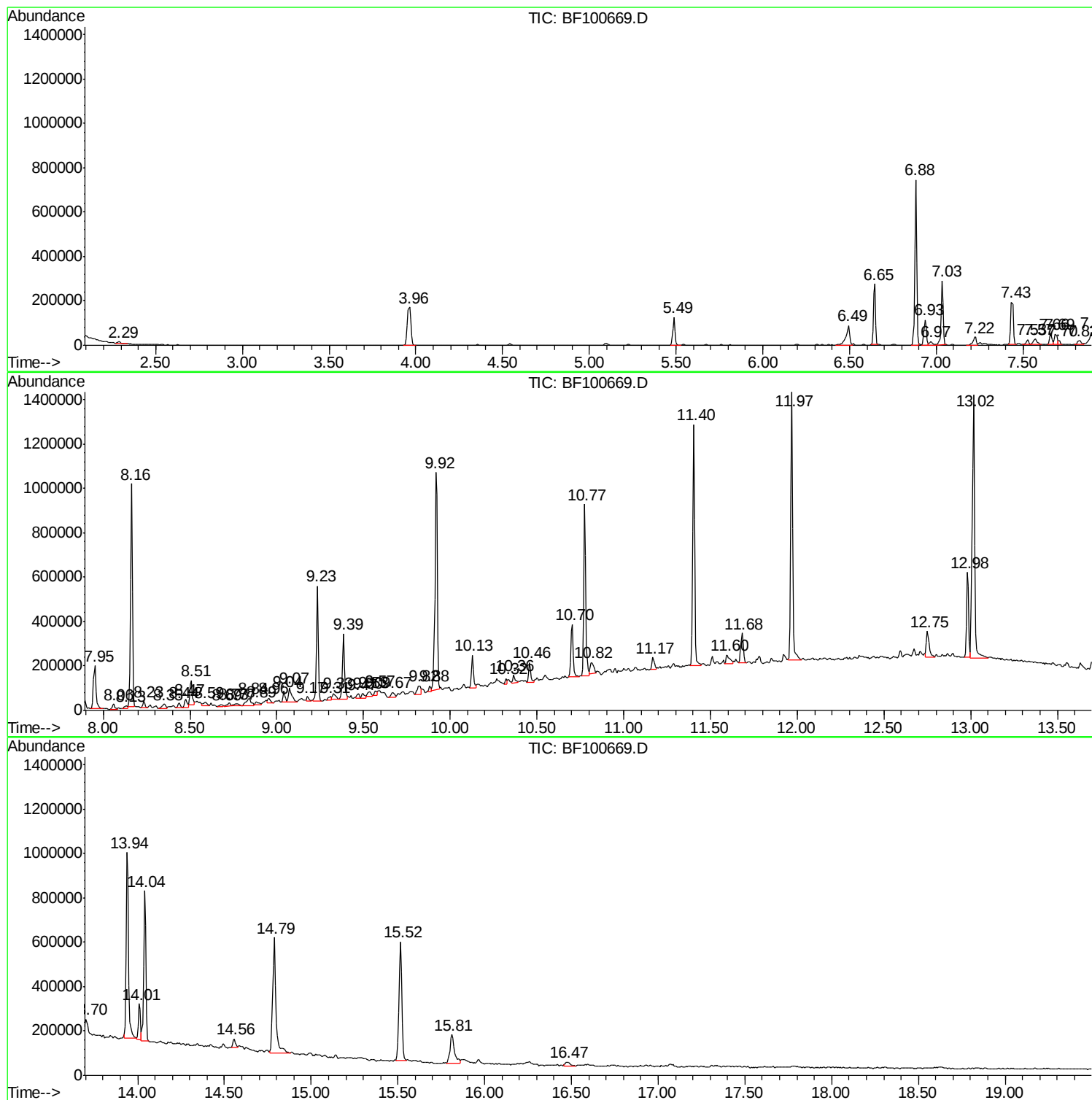
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Instrument :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF110817.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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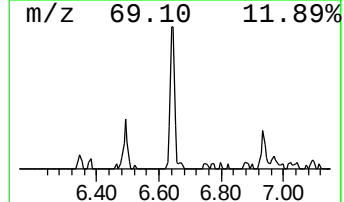
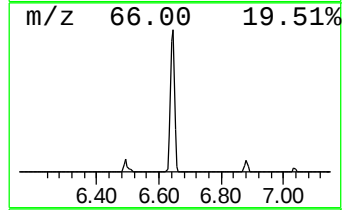
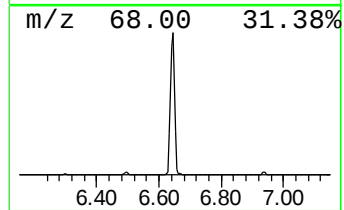
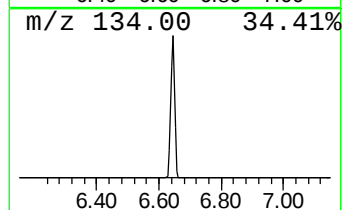
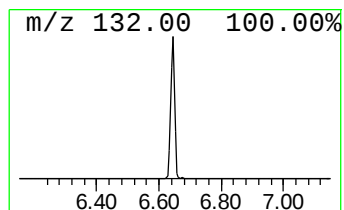
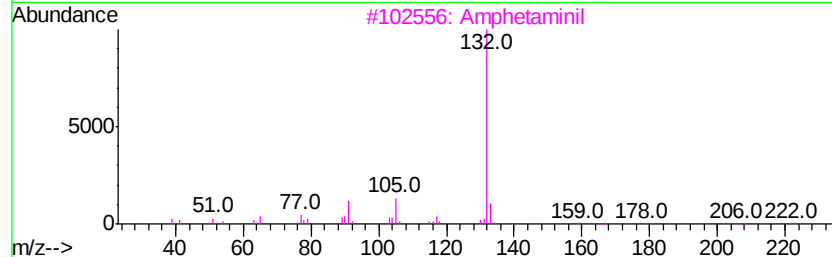
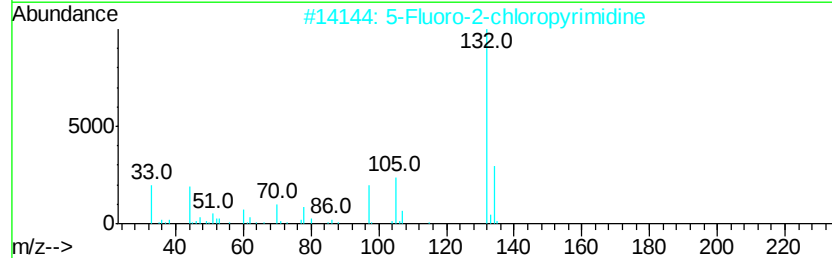
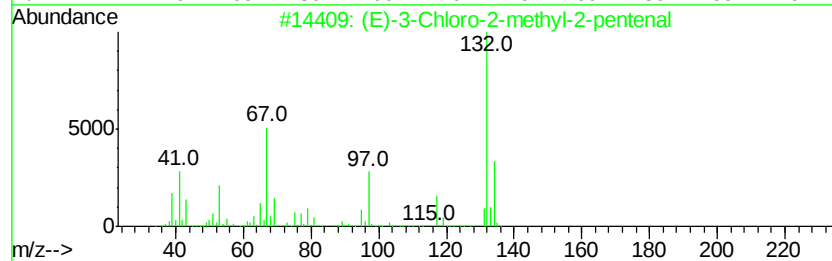
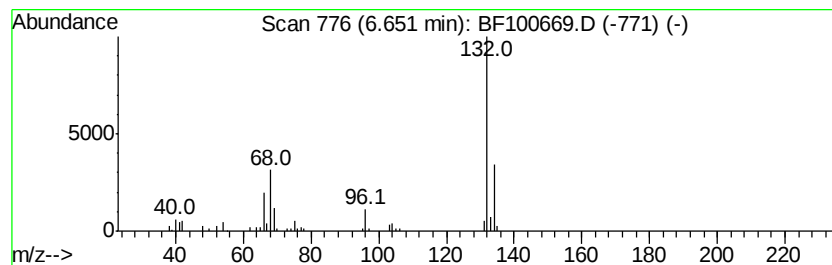
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF110817.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown6.65 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.65	7.81 ng	233722	1,4-Dichlorobenzene-d4	6.88

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	35
2			5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	17
3			Amphetaminil	250	C17H18N2	017590-01-1	12
4			Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
5			4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	9



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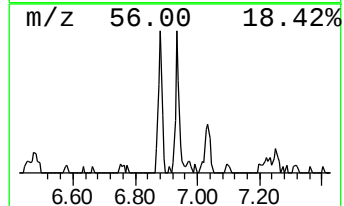
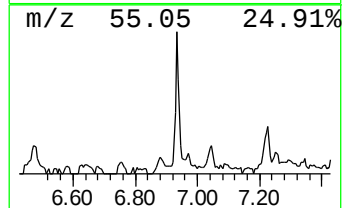
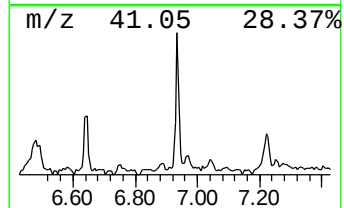
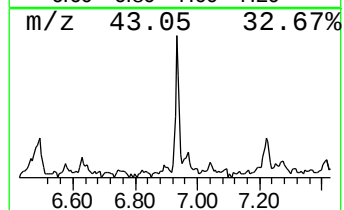
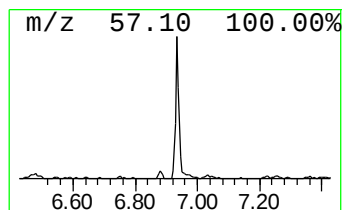
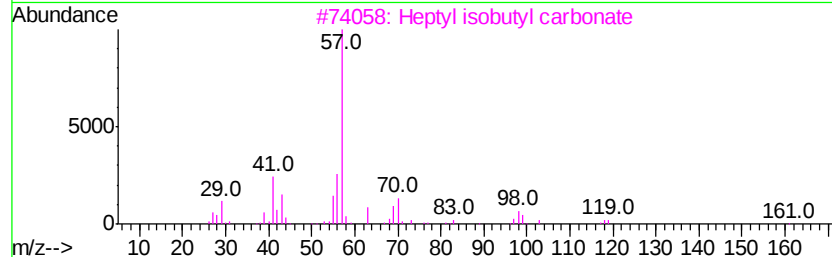
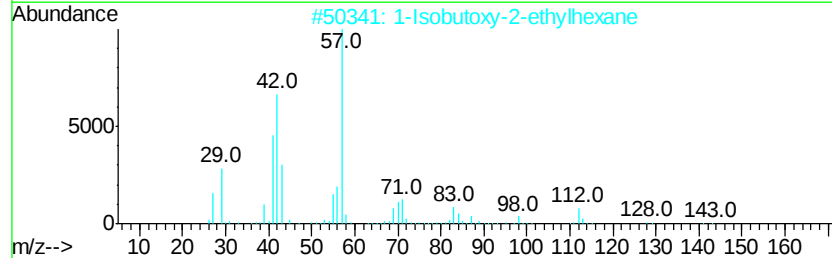
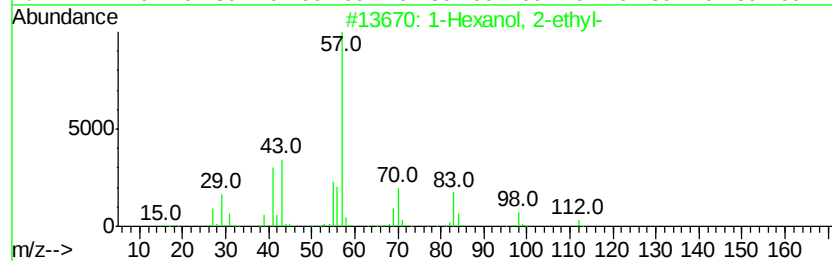
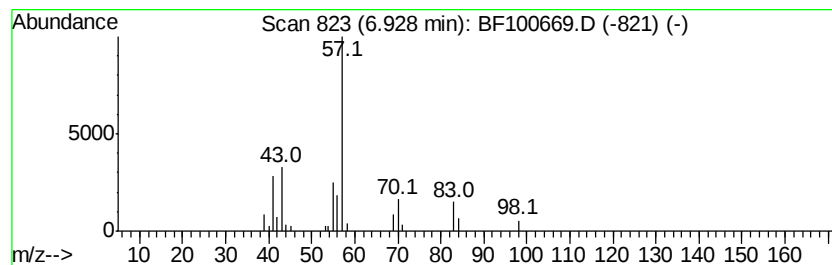
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF110817.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 1-Hexanol, 2-ethyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.93	3.12 ng	93297	1,4-Dichlorobenzene-d4	6.88

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	78
2			1-Isobutoxy-2-ethylhexane	186	C12H26O	1000139-90-3	64
3			Heptyl isobutyl carbonate	216	C12H24O3	959068-08-7	59
4			1-Pentanol, 2-ethyl-4-methyl-	130	C8H18O	000106-67-2	53
5			Heptane, 2,3,6-trimethyl-	142	C10H22	004032-93-3	53



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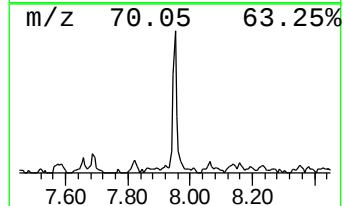
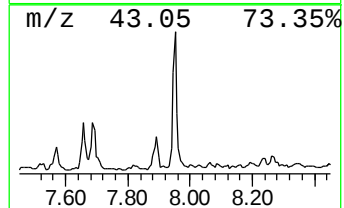
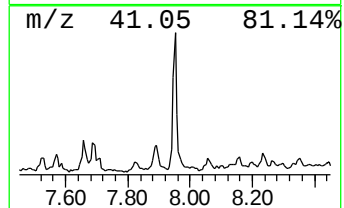
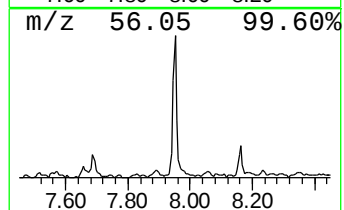
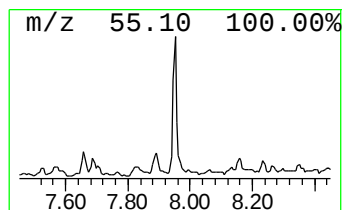
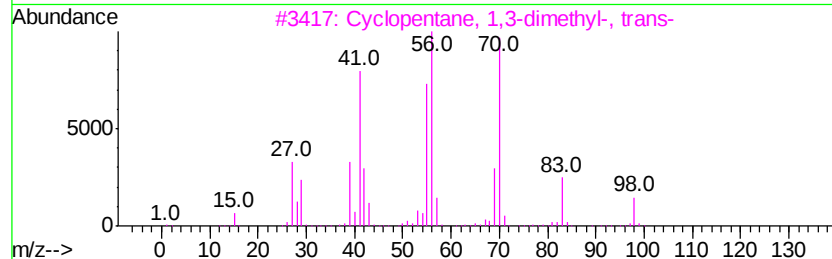
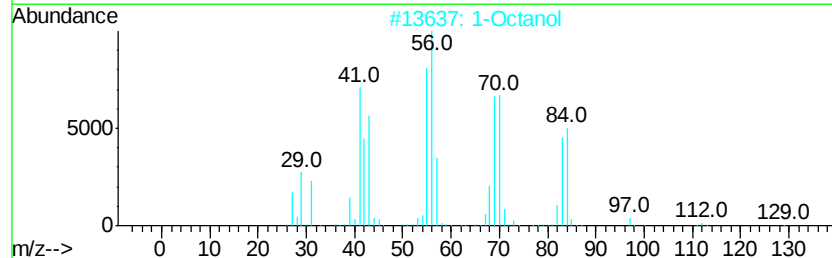
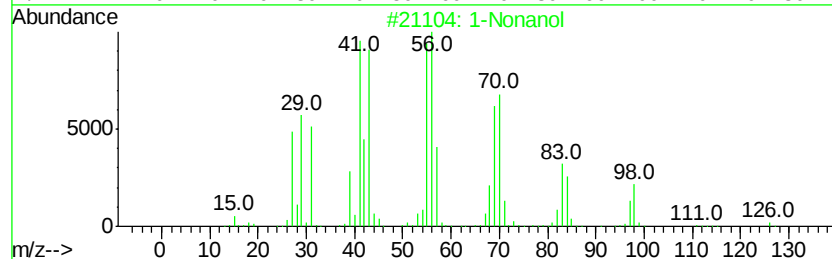
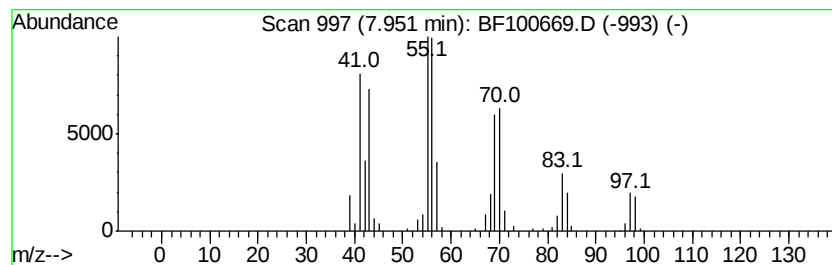
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 1-Nonanol Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.95	4.32 ng	176629	Napthalene-d8	8.16

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Nonanol	144	C9H20O	000143-08-8	91
2		1-Octanol	130	C8H18O	000111-87-5	80
3		Cyclopentane, 1,3-dimethyl-, trans-	98	C7H14	001759-58-6	53
4		Cyclopentane, 1,2-dimethyl-	98	C7H14	002452-99-5	52
5		Isopropylcyclobutane	98	C7H14	000872-56-0	52



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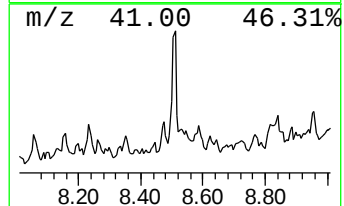
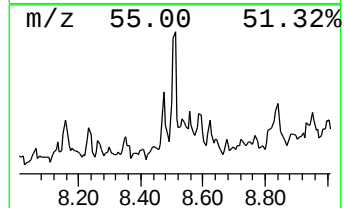
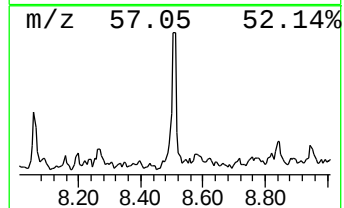
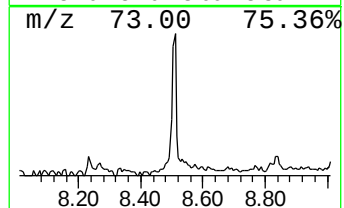
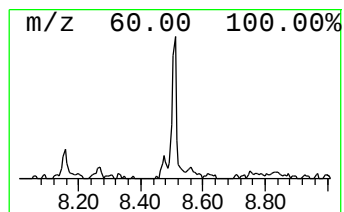
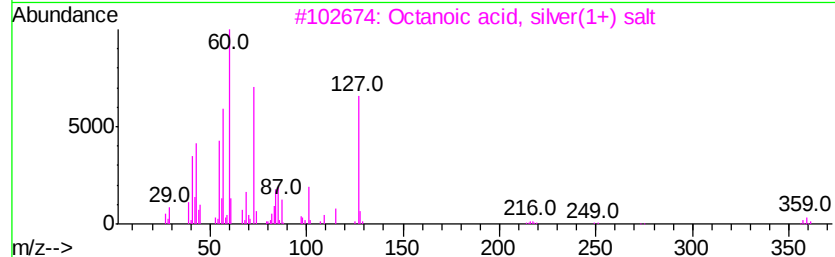
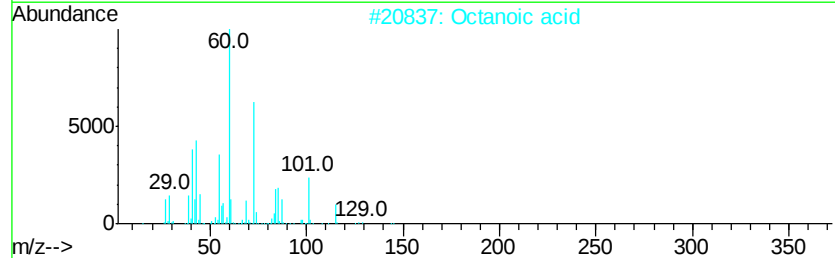
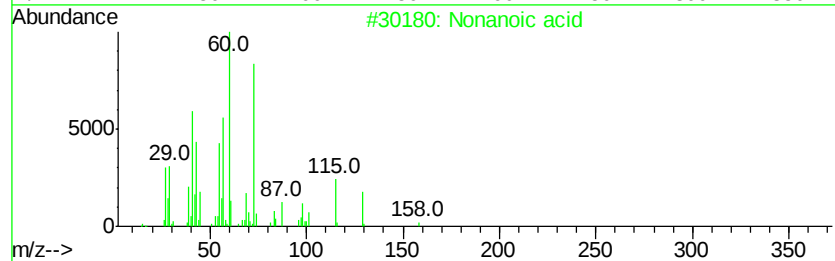
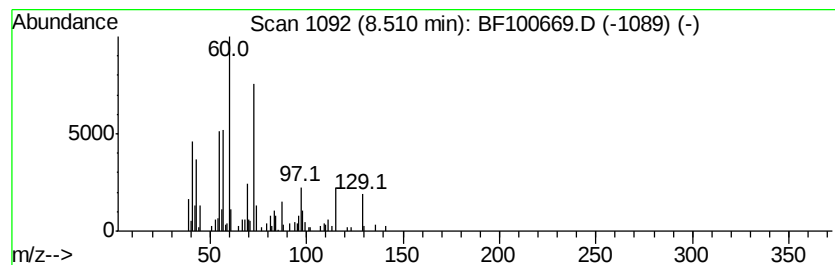
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Nonanoic acid Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.51	2.38 ng	97302	Napthalene-d8	8.16

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Nonanoic acid	158	C9H18O2	000112-05-0	72
2		Octanoic acid	144	C8H16O2	000124-07-2	50
3		Octanoic acid, silver(1+) salt	250	C8H15AgO2	024927-67-1	50
4		Pentanoic acid	102	C5H10O2	000109-52-4	47
5		Hexanoic acid	116	C6H12O2	000142-62-1	47



Data Path : Z:\HPCHEM1\BNA F\DATA\BF111717\
Data File : BF100669.D
Acq On : 17 Nov 2017 11:41
Operator : SJ/JU
Sample : I6477-01 10X
Misc :
ALS Vial : 6 Sample Multiplier: 1

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ClientSampleId :
TOTE-1-2

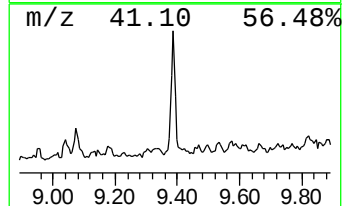
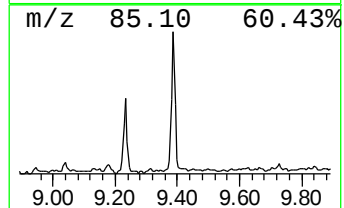
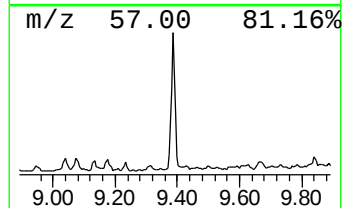
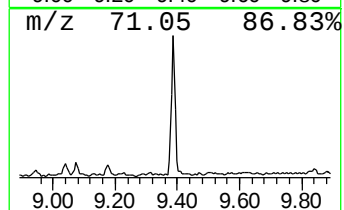
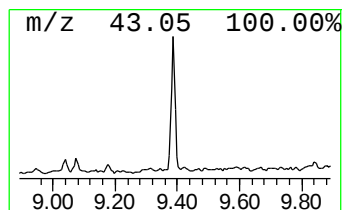
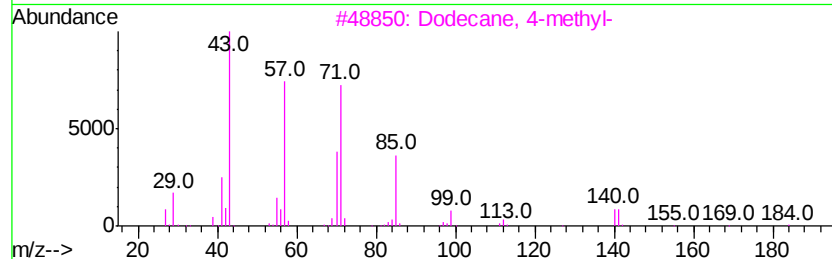
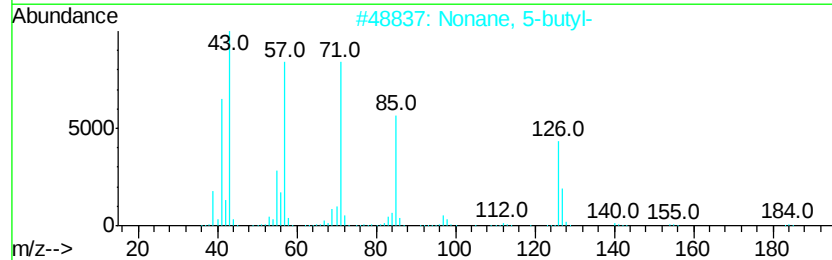
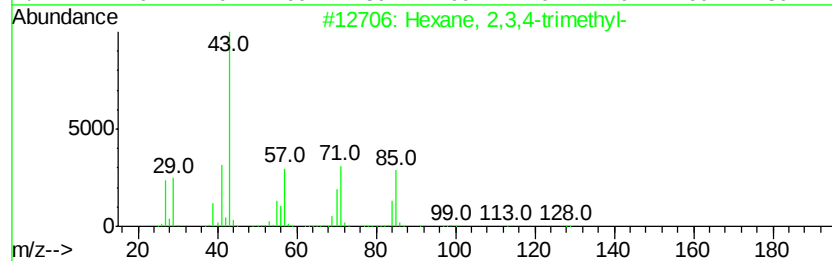
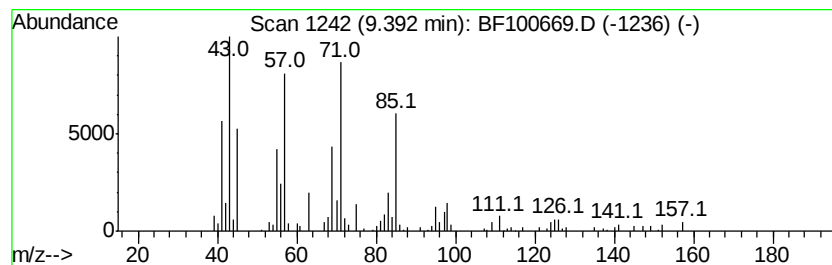
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Hexane, 2,3,4-trimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.39	5.70 ng	262206	Acenaphthene-d10	9.92

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexane, 2,3,4-trimethyl-	128	C9H20	000921-47-1	53
2		Nonane, 5-butyl-	184	C13H28	017312-63-9	53
3		Dodecane, 4-methyl-	184	C13H28	006117-97-1	47
4		Furan, tetrahydro-2-(methoxymeth...	116	C6H12O2	019354-27-9	43
5		Butane, 2,2-dimethyl-	86	C6H14	000075-83-2	38



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF111717\
Data File : BF100669.D
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ClientSampleId :
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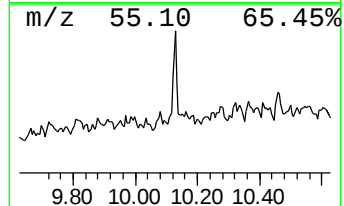
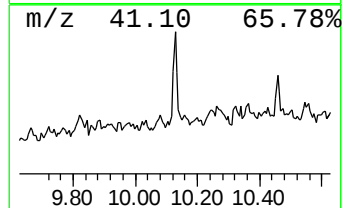
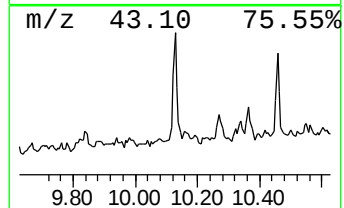
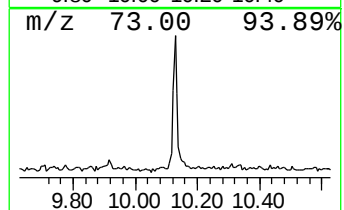
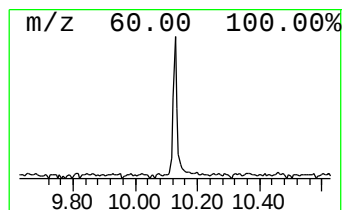
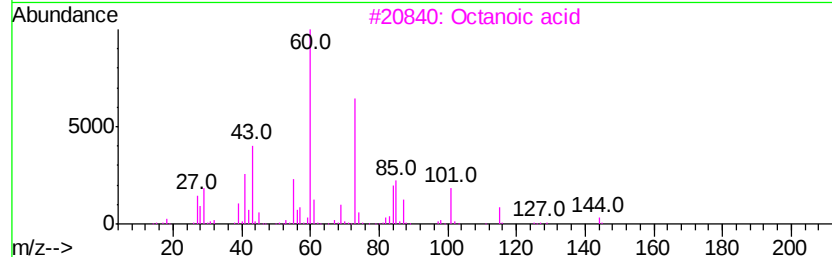
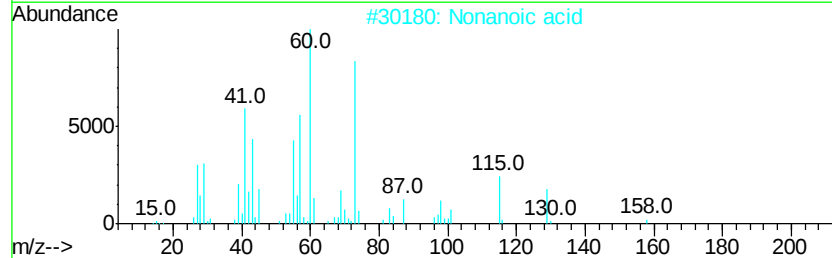
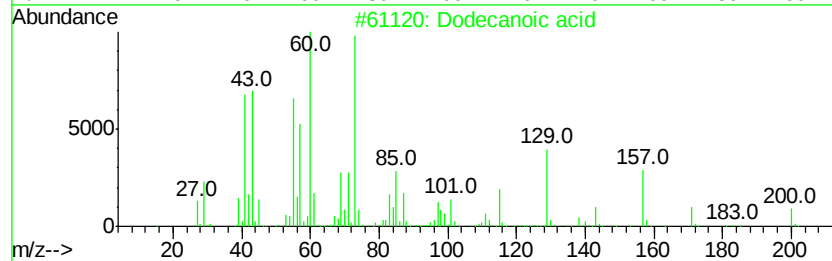
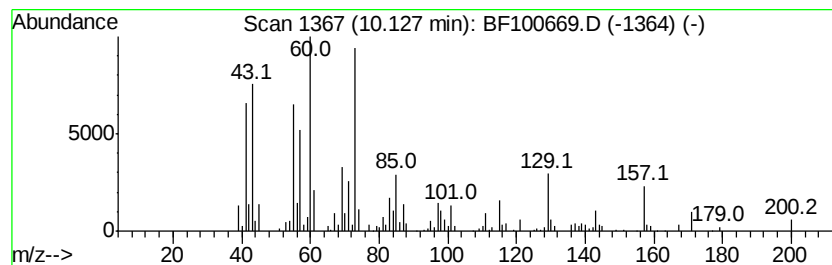
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Dodecanoic acid Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.13	2.87 ng	131838	Acenaphthene-d10	9.92

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Dodecanoic acid	200	C12H24O2	000143-07-7	97
2		Nonanoic acid	158	C9H18O2	000112-05-0	53
3		Octanoic acid	144	C8H16O2	000124-07-2	47
4		Ethanone, 1-(4,5-dihydro-2-thiaz...	129	C5H7NOS	029926-41-8	38
5		d-Glucohexodialdose	178	C6H10O6	1000149-19-1	38



Data Path : Z:\HPCHEM1\BNA F\DATA\BF111717\
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 Acq On : 17 Nov 2017 11:41
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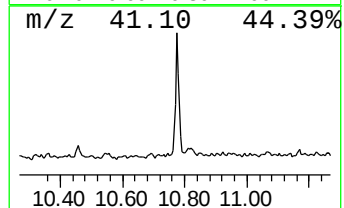
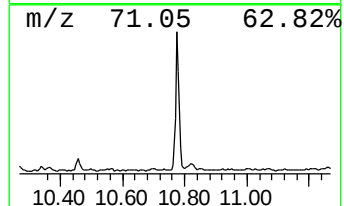
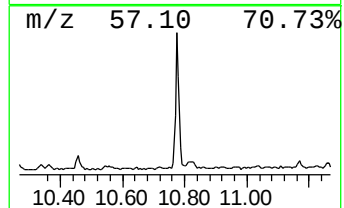
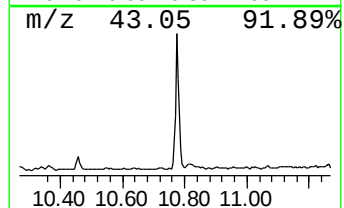
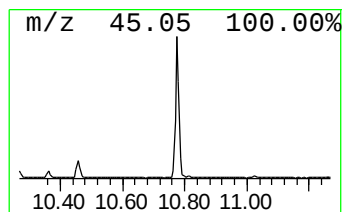
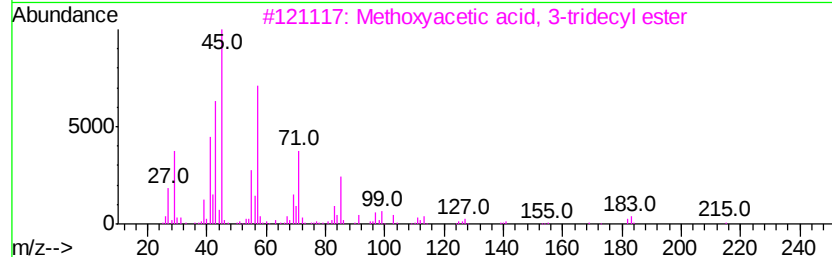
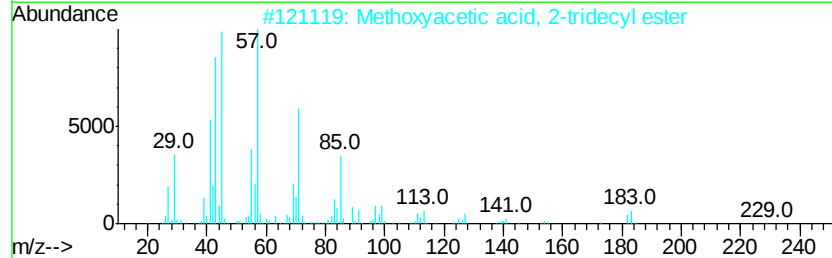
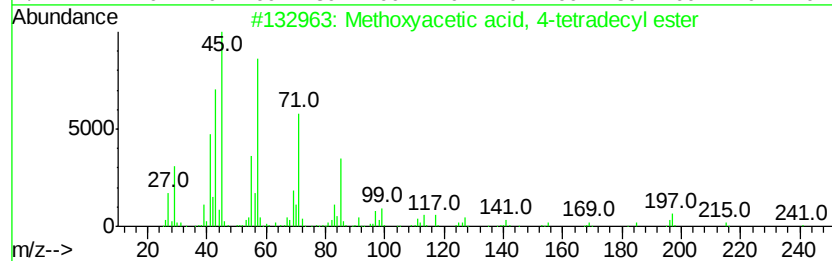
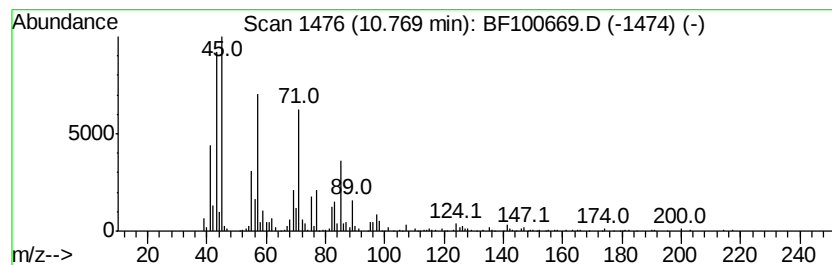
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 9 Methoxyacetic acid, 4-tetra... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.77	13.65 ng	635560	Phenanthrene-d10	11.40

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Methoxvacetic acid, 4-tetradecyl...	286	C17H34O3	1000282-05-0	72
2		Methoxvacetic acid, 2-tridecyl e...	272	C16H32O3	1000282-04-5	58
3		Methoxvacetic acid, 3-tridecyl e...	272	C16H32O3	1000282-04-6	47
4		3-Ethyl-3-methyl-2-pentanol	130	C8H18O	066576-22-5	43
5		Diethylene glycol hexyl ether	190	C10H22O3	000112-59-4	38



Data Path : Z:\HPCHEM1\BNA F\DATA\BF111717\
Data File : BF100669.D
Acq On : 17 Nov 2017 11:41
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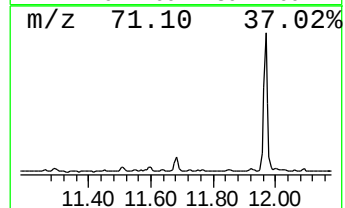
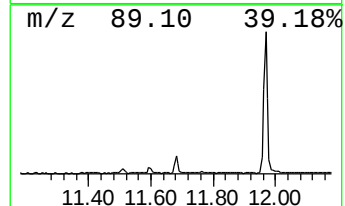
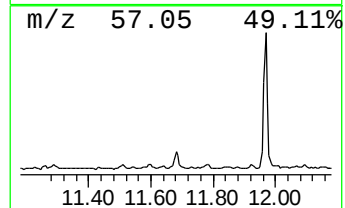
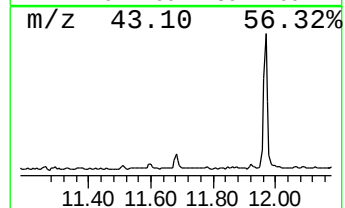
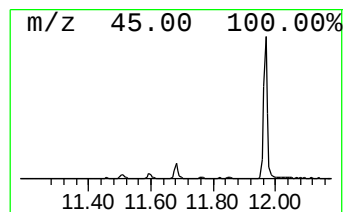
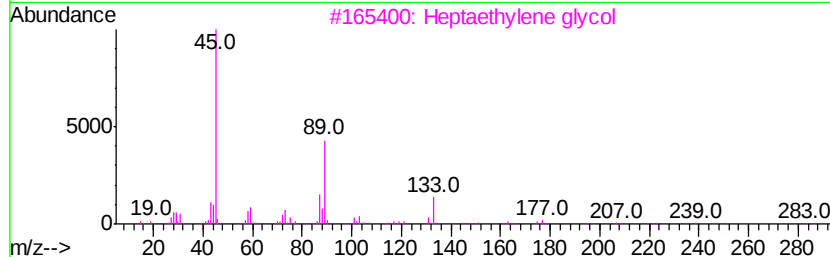
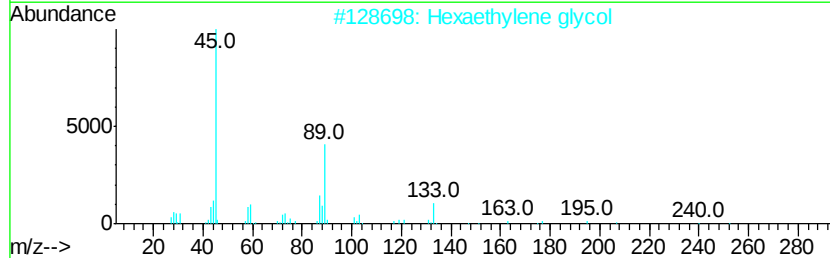
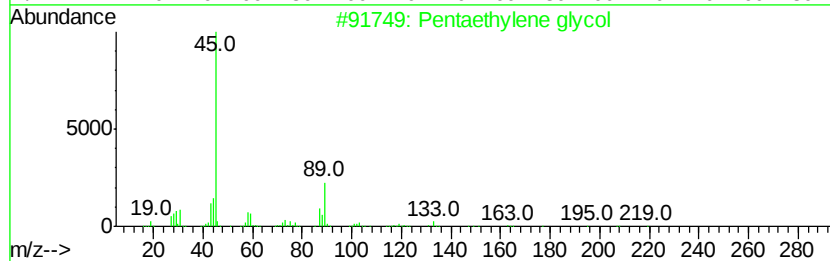
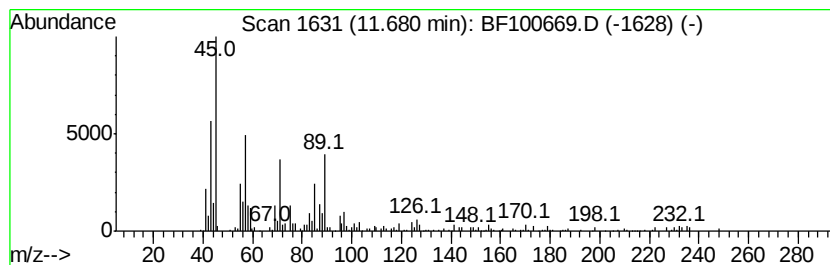
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Pentaethylene glycol Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.68	2.49 ng	115786	Phenanthrene-d10	11.40

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentaethylene glycol	238	C10H22O6	004792-15-8	58
2			Hexaethylene glycol	282	C12H26O7	002615-15-8	58
3			Heptaethylene glycol	326	C14H30O8	005617-32-3	53
4			2-[2-[2-[2-[2-[2-(2-Hydroxyet...	370	C16H34O9	005117-19-1	53
5			Tetraethylene glycol	194	C8H18O5	000112-60-7	53



Data Path : Z:\HPCHEM1\BNA F\DATA\BF111717\
Data File : BF100669.D
Acq On : 17 Nov 2017 11:41
Operator : SJ/JU
Sample : I6477-01 10X
Misc :
ALS Vial : 6 Sample Multiplier: 1

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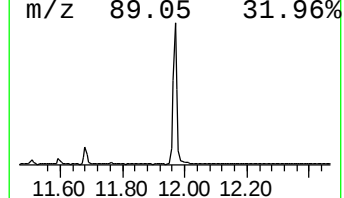
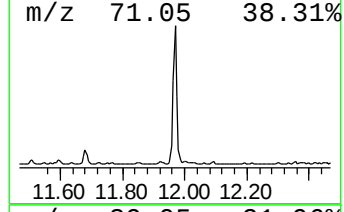
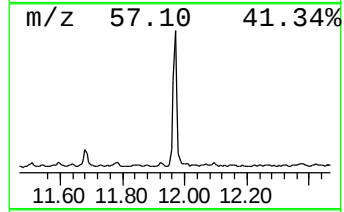
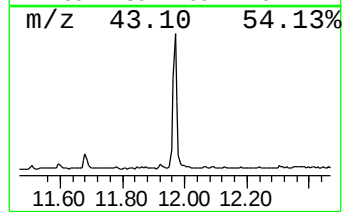
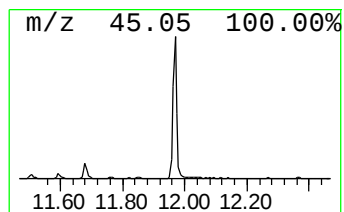
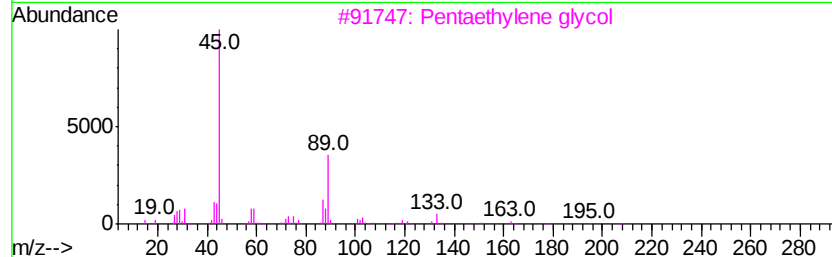
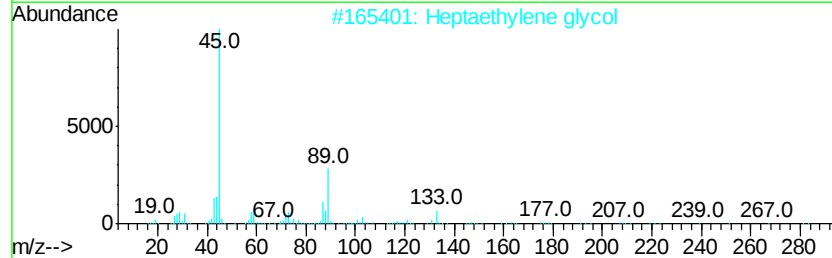
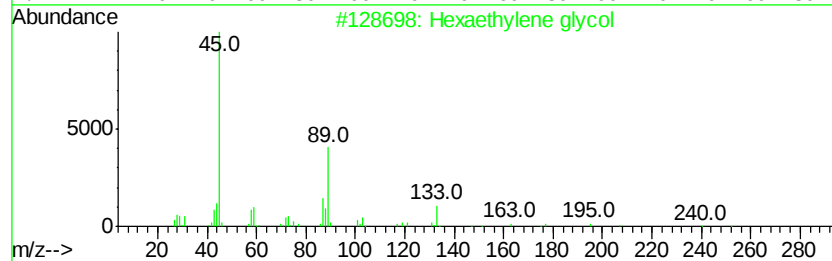
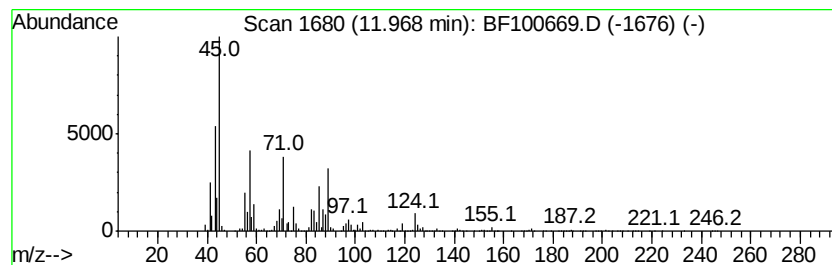
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 unknown11.97 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.97	22.29 ng	1038320	Phenanthrene-d10	11.40

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexaethylene glycol	282	C12H26O7	002615-15-8	49
2		Heptaethylene glycol	326	C14H30O8	005617-32-3	49
3		Pentaethylene glycol	238	C10H22O6	004792-15-8	47
4		Ethanol, 2-[2-(2-methoxyethoxy)ethoxy]-	164	C7H16O4	000112-35-6	43
5		Ethanol, 2-[2-(ethenyloxy)ethoxy]-	132	C6H12O3	000929-37-3	38



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF111717\
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ALS Vial : 6 Sample Multiplier: 1

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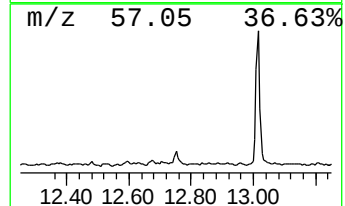
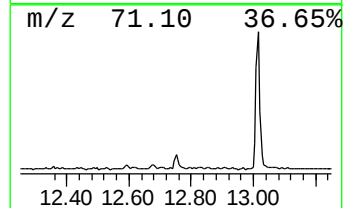
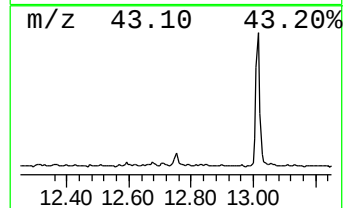
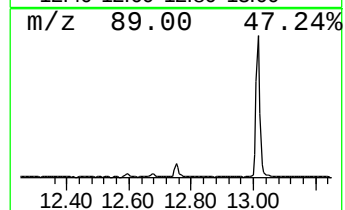
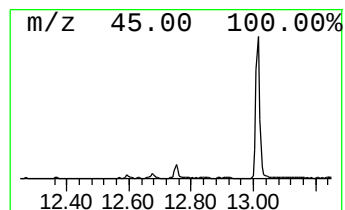
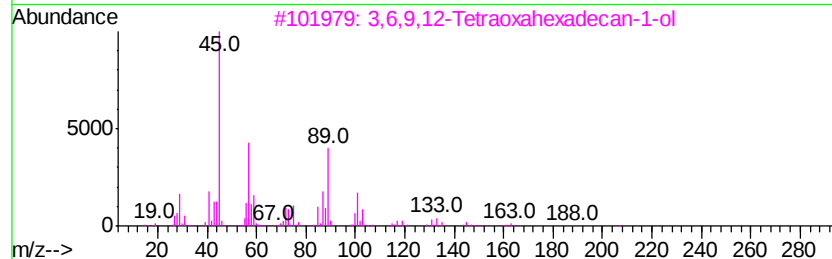
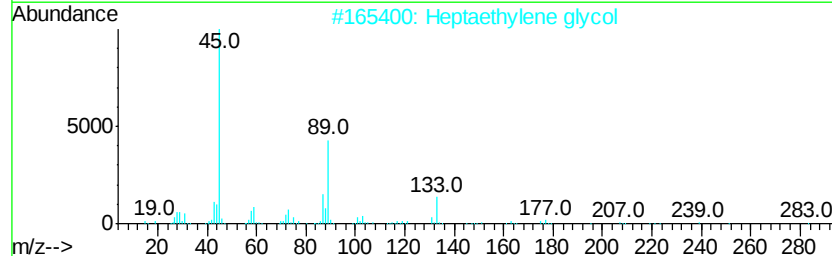
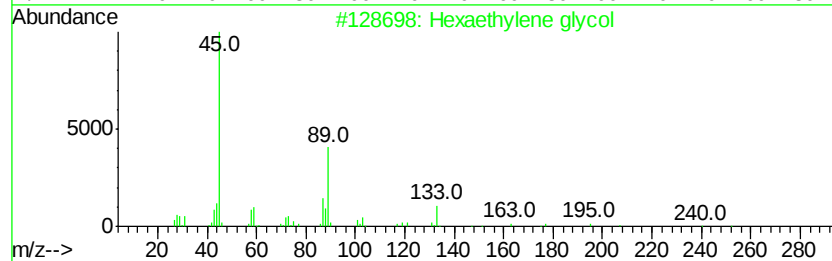
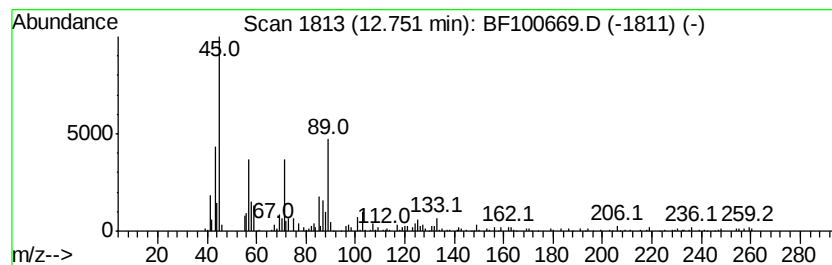
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Hexaethylene glycol Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.75	3.43 ng	105084	Chrysene-d12	14.04

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexaethylene glycol	282	C12H26O7	002615-15-8	72
2		Heptaethylene glycol	326	C14H30O8	005617-32-3	72
3		3,6,9,12-Tetraoxahexadecan-1-ol	250	C12H26O5	001559-34-8	72
4		Pentaethylene glycol	238	C10H22O6	004792-15-8	64
5		Pentaethylene glycol monododecyl...	406	C22H46O6	003055-95-6	64



Data Path : Z:\HPCHEM1\BNA F\DATA\BF111717\
Data File : BF100669.D
Acq On : 17 Nov 2017 11:41
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Sample : I6477-01 10X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
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ClientSampleId :
TOTE-1-2

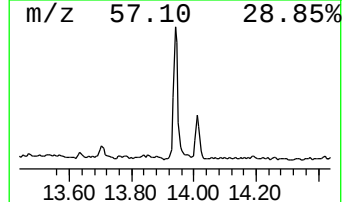
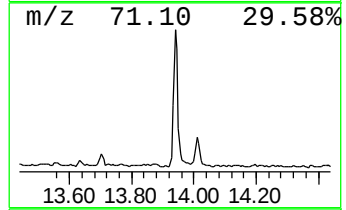
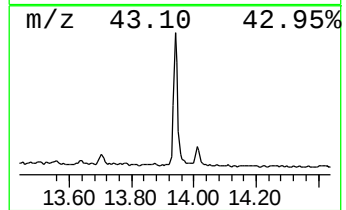
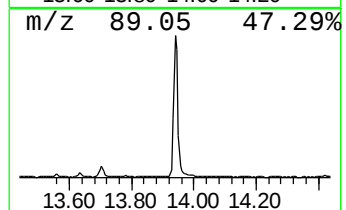
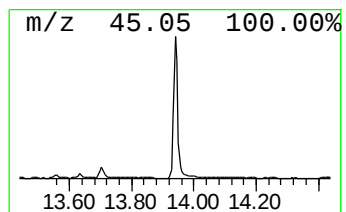
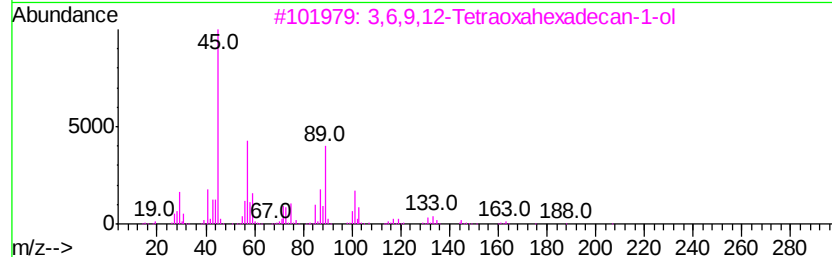
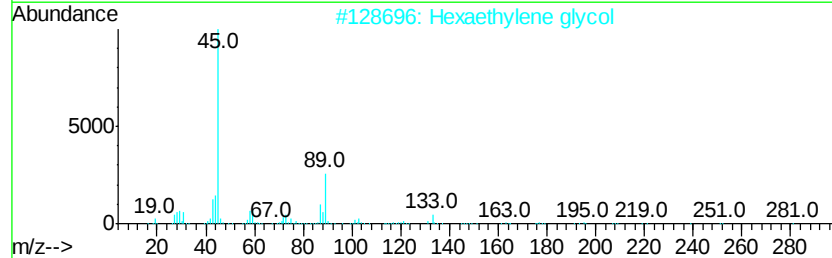
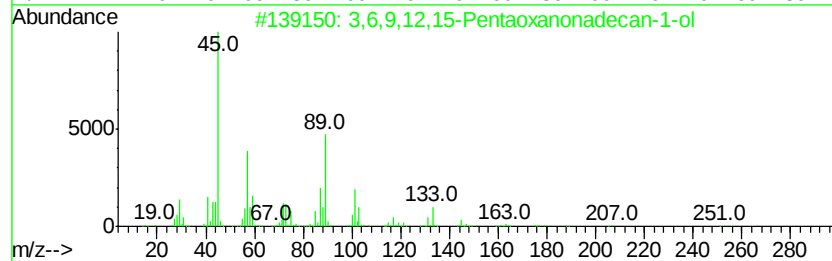
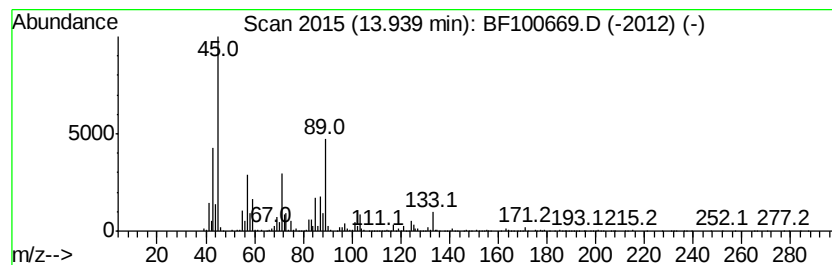
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 3,6,9,12,15-Pentaoxanonadec... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.94	26.88 ng	823694	Chrysene-d12	14.04

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3,6,9,12,15-Pentaoxanonadecan-1-ol	294	C14H30O6	001786-94-3	83
2		Hexaethylene glycol	282	C12H26O7	002615-15-8	72
3		3,6,9,12-Tetraoxahexadecan-1-ol	250	C12H26O5	001559-34-8	72
4		Heptaethylene glycol	326	C14H30O8	005617-32-3	72
5		Hexaethylene glycol monododecyl ...	450	C24H50O7	003055-96-7	72



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF111717\
Data File : BF100669.D
Acq On : 17 Nov 2017 11:41
Operator : SJ/JU
Sample : I6477-01 10X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TOTE-1-2

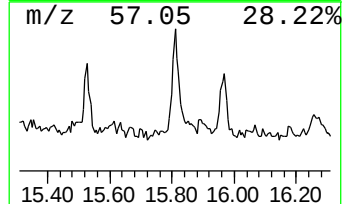
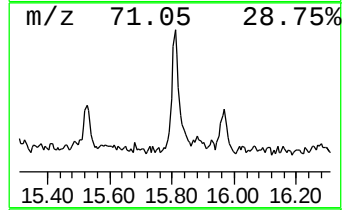
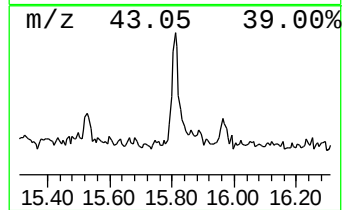
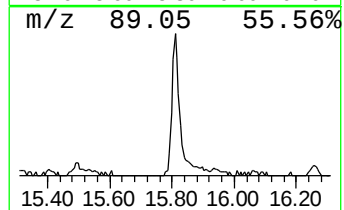
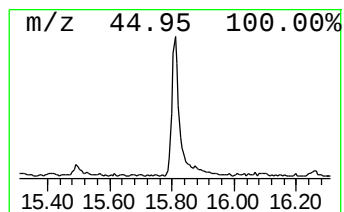
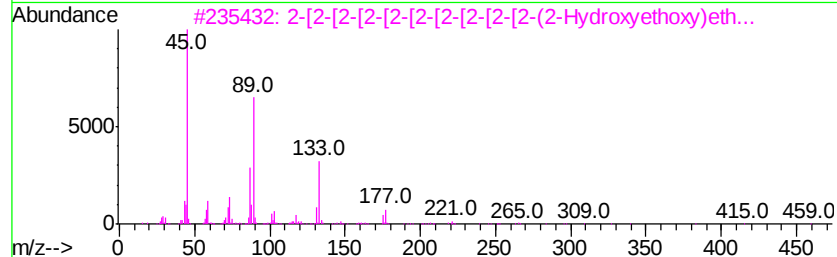
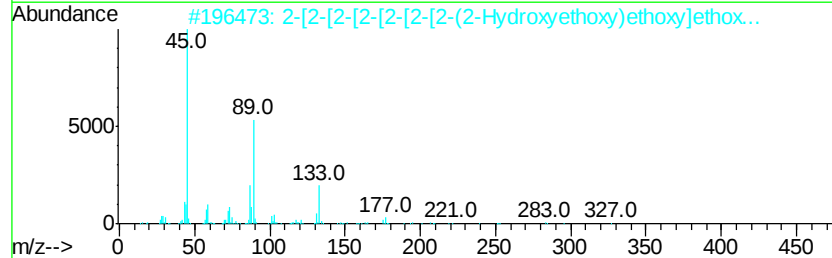
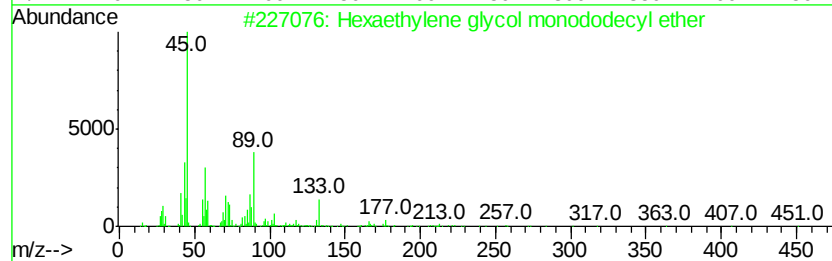
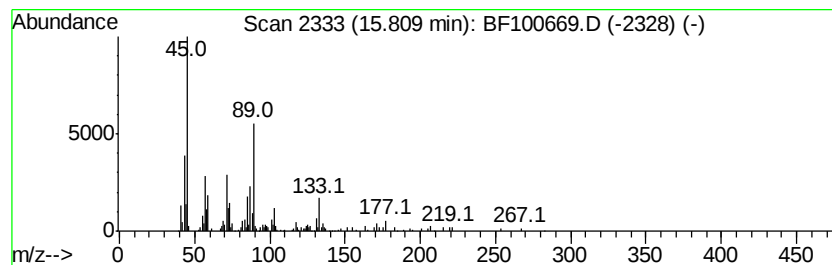
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 Hexaethylene glycol monododecyl... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.81	6.92 ng	228811	Perylene-d12	15.52

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexaethylene glycol monododecyl ...	450	C24H50O7	003055-96-7	87
2			2-[2-[2-[2-[2-[2-(2-Hydroxyet...	370	C16H34O9	005117-19-1	87
3			2-[2-[2-[2-[2-[2-[2-[2-(2-...	502	C22H46O12	1000352-13-2	83
4			3,6,9,12,15-Pentaoxanonadecan-1-ol	294	C14H30O6	001786-94-3	83
5			3,6,9,12-Tetraoxahexadecan-1-ol	250	C12H26O5	001559-34-8	80



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF111717\
Data File : BF100669.D
Acq On : 17 Nov 2017 11:41
Operator : SJ/JU
Sample : I6477-01 10X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TOTE-1-2

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF110817.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown6.65	6.65	7.8 ng		233722	1	6.88	598821	20.0
1-Hexanol, 2-ethyl-	6.93	3.1 ng		93297	1	6.88	598821	20.0
1-Nonanol	7.95	4.3 ng		176629	2	8.16	818379	20.0
Nonanoic acid	8.51	2.4 ng		97302	2	8.16	818379	20.0
Hexane, 2,3,4-tri...	9.39	5.7 ng		262206	3	9.92	920027	20.0
Dodecanoic acid	10.13	2.9 ng		131838	3	9.92	920027	20.0
Methoxyacetic aci...	10.77	13.7 ng		635560	4	11.40	931472	20.0
Pentaethylene glycol	11.68	2.5 ng		115786	4	11.40	931472	20.0
unknown11.97	11.97	22.3 ng		1038320	4	11.40	931472	20.0
Hexaethylene glycol	12.75	3.4 ng		105084	5	14.04	612916	20.0
3,6,9,12,15-Penta...	13.94	26.9 ng		823694	5	14.04	612916	20.0
Hexaethylene glyc...	15.81	6.9 ng		228811	6	15.52	660979	20.0