

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF052615\
 Data File : BF079507.D
 Acq On : 27 May 2015 7:08
 Operator : TP/IZ
 Sample : G2359-08
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-30AS

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-BF052615.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	1.278	5	8	10	rVB	86207	128135	5.95%	0.512%
2	1.427	19	21	24	rVB	201114	231896	10.78%	0.927%
3	1.518	27	29	34	rVB	25561	34007	1.58%	0.136%
4	1.587	34	35	43	rVV	49321	112509	5.23%	0.450%
5	1.701	43	45	68	rVB	590111	1149701	53.43%	4.594%
6	4.742	309	311	321	rVB	114659	161380	7.50%	0.645%
7	5.302	358	360	373	rBV	1189444	1411961	65.62%	5.642%
8	6.616	472	475	483	rBV	1436609	1536844	71.42%	6.141%
9	6.730	483	485	488	rBV	1176455	1513240	70.33%	6.046%
10	7.016	508	510	513	rBV	440686	437910	20.35%	1.750%
11	7.199	524	526	529	rBV	968770	1143005	53.12%	4.567%
12	7.713	569	571	574	rBV	771422	867233	40.30%	3.465%
13	8.593	646	648	651	rBV	634172	616413	28.65%	2.463%
14	9.942	764	766	769	rBV	1969232	1810405	84.14%	7.234%
15	10.456	808	811	813	rBV	71996	70302	3.27%	0.281%
16	10.594	820	823	825	rBV	21484	27579	1.28%	0.110%
17	10.731	831	835	836	rBV2	10274	24048	1.12%	0.096%
18	10.765	836	838	840	rVV	745697	723572	33.63%	2.891%
19	10.799	840	841	846	rVB3	26458	28189	1.31%	0.113%
20	11.359	888	890	892	rVB	23619	21794	1.01%	0.087%
21	11.439	896	897	901	rBV	16483	24061	1.12%	0.096%
22	11.748	921	924	926	rBV	1167371	1316631	61.19%	5.261%
23	11.942	939	941	945	rBV2	14920	24414	1.13%	0.098%
24	12.605	996	999	1000	rBV	621535	789772	36.70%	3.156%
25	12.628	1000	1001	1003	rVV	298646	255250	11.86%	1.020%
26	12.697	1003	1007	1009	rVB	72264	88657	4.12%	0.354%
27	13.234	1051	1054	1055	rBV	32217	40689	1.89%	0.163%
28	13.268	1055	1057	1060	rVB	37953	54819	2.55%	0.219%
29	13.360	1060	1065	1067	rBV2	101751	218301	10.15%	0.872%
30	13.394	1067	1068	1071	rVB	32704	29546	1.37%	0.118%
31	13.600	1085	1086	1088	rBV	20607	28763	1.34%	0.115%
32	13.920	1111	1114	1115	rBV	25283	37596	1.75%	0.150%
33	13.954	1115	1117	1121	rVB2	19778	44728	2.08%	0.179%
34	14.011	1121	1122	1125	rVB2	23496	42498	1.98%	0.170%

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Integrator: RTE

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Sampling : 1

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If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF052615.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	14.103	1128	1130	1132	rBV	541248	522551	24.28%	2.088%
36	14.285	1143	1146	1148	rBV	30259	33299	1.55%	0.133%
37	14.377	1151	1154	1157	rBV	484407	583330	27.11%	2.331%
38	14.548	1166	1169	1170	rBV	24217	34152	1.59%	0.136%
39	14.594	1170	1173	1176	rVV	2024345	2151772	100.00%	8.598%
40	14.674	1176	1180	1182	rVB2	24809	47918	2.23%	0.191%
41	14.765	1185	1188	1189	rBV3	25895	43404	2.02%	0.173%
42	14.800	1189	1191	1194	rVB	63993	83125	3.86%	0.332%
43	14.880	1196	1198	1200	rBV	34933	60054	2.79%	0.240%
44	14.925	1200	1202	1207	rVB2	61662	115999	5.39%	0.463%
45	15.040	1210	1212	1213	rBV	25055	23471	1.09%	0.094%
46	15.074	1213	1215	1218	rBV	30461	46800	2.17%	0.187%
47	15.246	1227	1230	1232	rBV	171579	214358	9.96%	0.857%
48	15.440	1243	1247	1250	rVB3	26586	55777	2.59%	0.223%
49	15.508	1252	1253	1255	rBV2	20365	23556	1.09%	0.094%
50	15.566	1255	1258	1260	rBV	46559	65078	3.02%	0.260%
51	15.600	1260	1261	1262	rBV	37811	39075	1.82%	0.156%
52	15.703	1268	1270	1273	rVB	44286	49166	2.28%	0.196%
53	15.771	1273	1276	1280	rBV2	272349	419577	19.50%	1.677%
54	15.874	1282	1285	1287	rVV2	761867	1258601	58.49%	5.029%
55	15.908	1287	1288	1293	rVB2	230093	360448	16.75%	1.440%
56	16.000	1294	1296	1298	rBV	39310	63483	2.95%	0.254%
57	16.046	1298	1300	1301	rBV	53866	68994	3.21%	0.276%
58	16.137	1307	1308	1310	rVB2	26071	27990	1.30%	0.112%
59	16.183	1310	1312	1316	rBV2	62497	74737	3.47%	0.299%
60	16.286	1319	1321	1323	rBV2	16640	33452	1.55%	0.134%
61	16.331	1323	1325	1326	rBV2	28151	38252	1.78%	0.153%
62	16.377	1326	1329	1331	rBV	46970	60288	2.80%	0.241%
63	16.411	1331	1332	1334	rBV	18021	29814	1.39%	0.119%
64	16.491	1337	1339	1340	rVV	29053	32201	1.50%	0.129%
65	16.526	1340	1342	1345	rVV3	27096	79326	3.69%	0.317%
66	16.571	1345	1346	1349	rVB2	98675	148074	6.88%	0.592%
67	16.914	1373	1376	1377	rBV3	29446	59967	2.79%	0.240%
68	16.960	1378	1380	1385	rVV2	51636	156882	7.29%	0.627%
69	17.154	1395	1397	1403	rVB	268392	536779	24.95%	2.145%
70	17.269	1405	1407	1409	rVB	59267	68792	3.20%	0.275%
71	17.337	1411	1413	1419	rBV3	75543	235345	10.94%	0.940%

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ClientSampleId :
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Integration Parameters: rteint.p
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

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

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Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

72	17.463	1422	1424	1427	rBV	165438	217481	10.11%	0.869%
73	17.531	1427	1430	1433	rVB	220301	279959	13.01%	1.119%
74	17.589	1433	1435	1437	rBV	744875	910085	42.29%	3.636%
75	18.914	1548	1551	1560	rVB2	118001	257530	11.97%	1.029%
76	19.303	1582	1585	1589	rVB	98946	188344	8.75%	0.753%
77	19.440	1594	1597	1602	rBV5	76031	181817	8.45%	0.726%

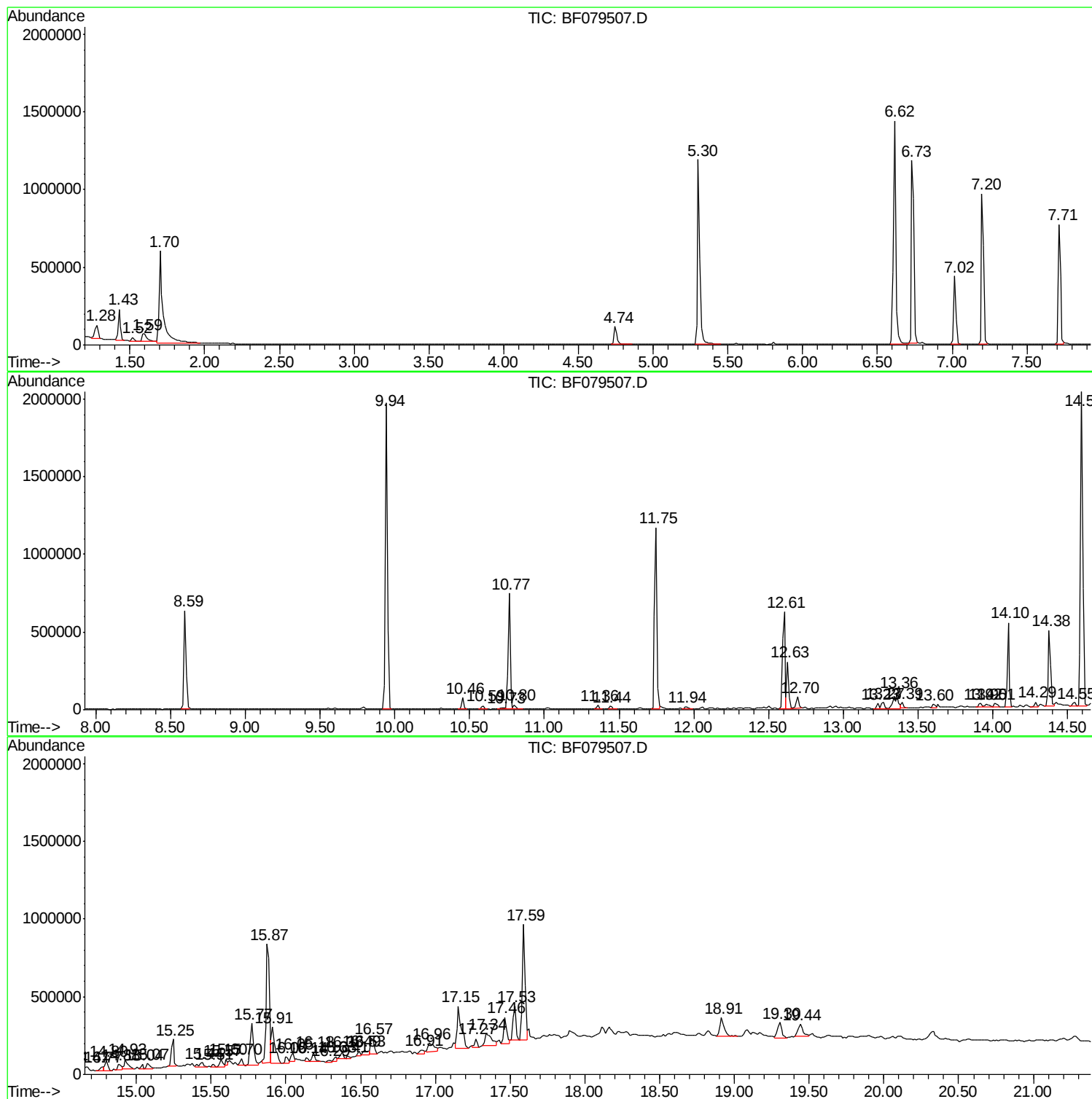
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Instrument :
BNA_F
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF052615.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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Acq On : 27 May 2015 7:08
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Instrument :
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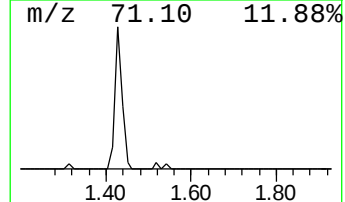
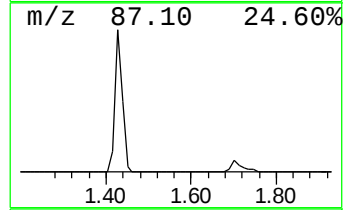
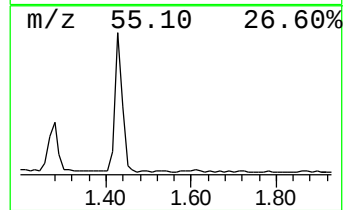
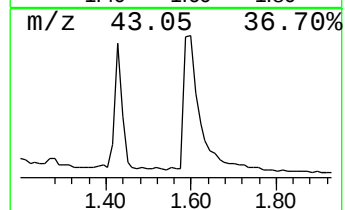
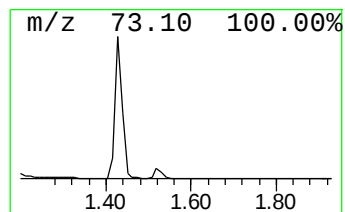
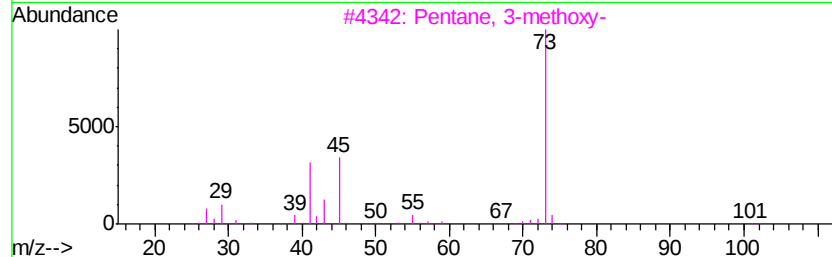
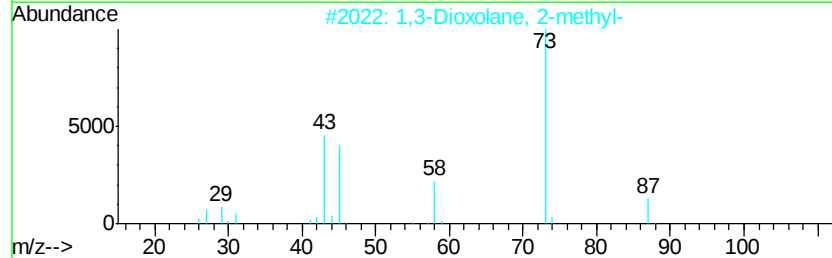
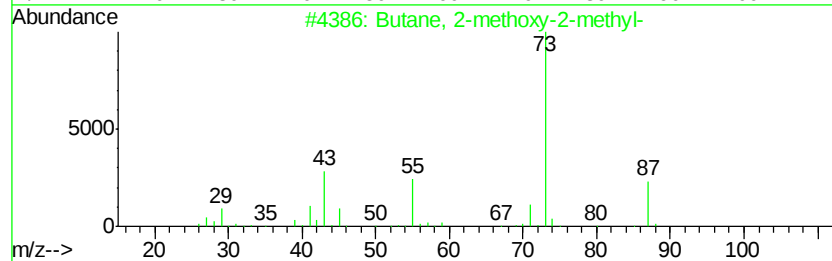
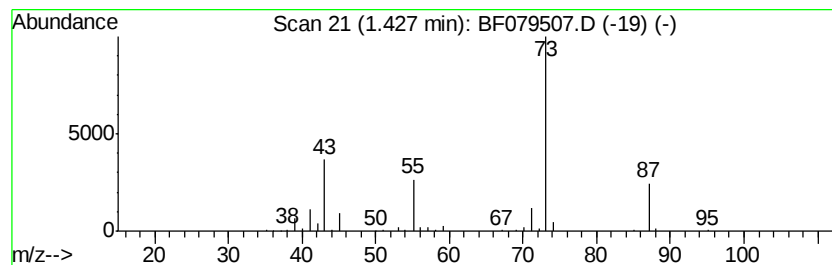
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF052615.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Butane, 2-methoxy-2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.43	10.59 ng	231896	1,4-Dichlorobenzene-d4	7.02

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2		1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	23
3		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	17
4		Silane, tetramethyl-	88	C4H12Si	000075-76-3	12
5		Borane, dimethoxy-	74	C2H7BO2	004542-61-4	9



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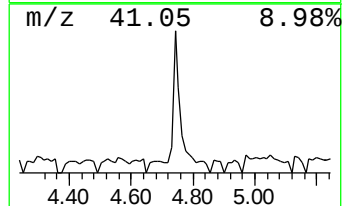
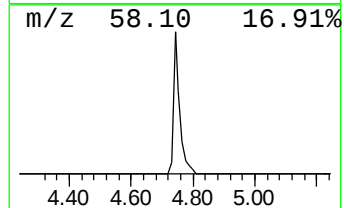
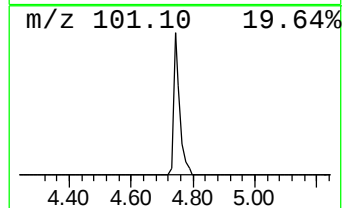
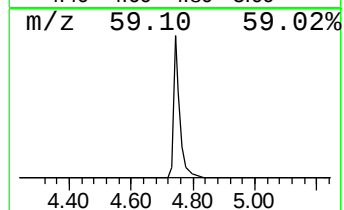
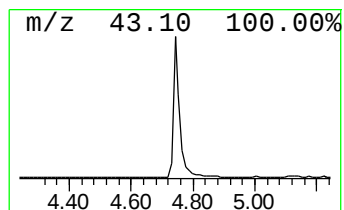
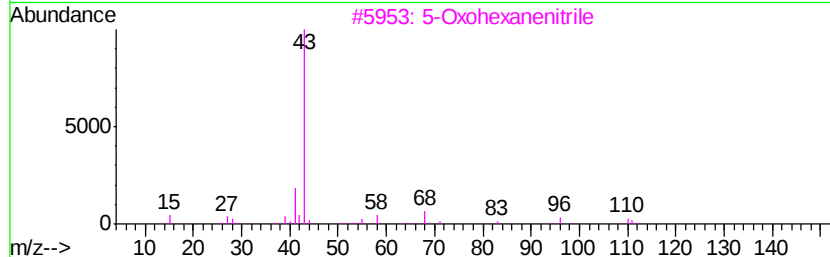
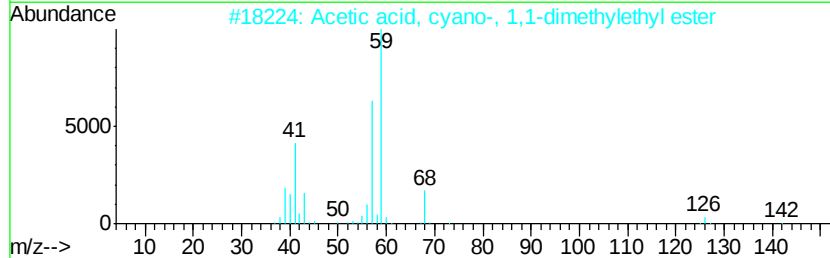
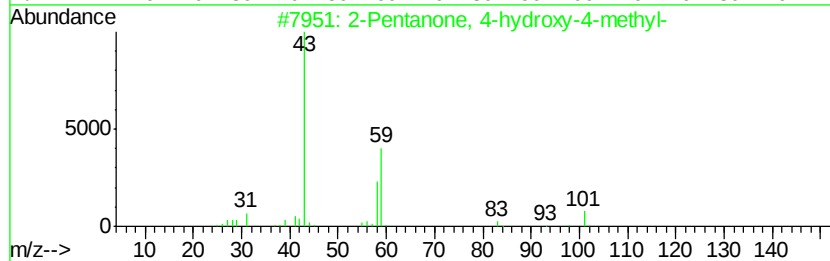
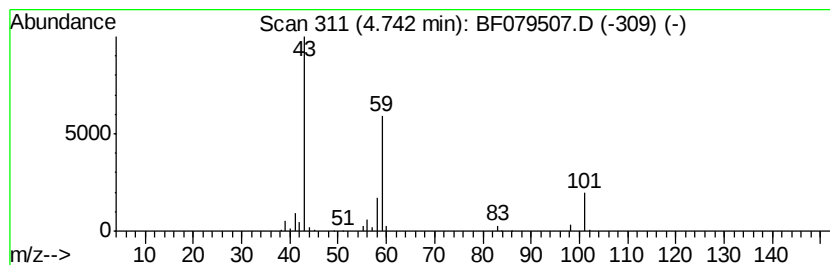
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF052615.M
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.74	7.37 ng	161380	1,4-Dichlorobenzene-d4	7.02

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2		Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	23
3		5-Oxohexanenitrile	111	C6H9NO	010412-98-3	9
4		t-Butyl ethyl ether	102	C6H14O	1000118-18-4	9
5		4-Hydroxy-3-hexanone	116	C6H12O2	004984-85-4	9



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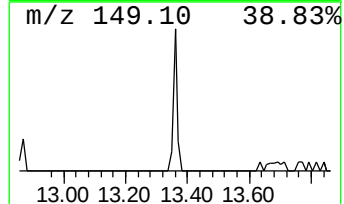
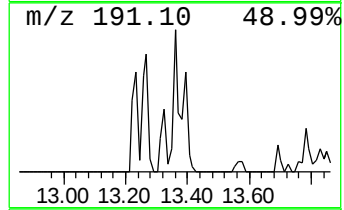
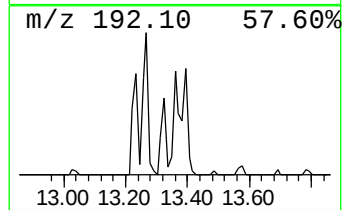
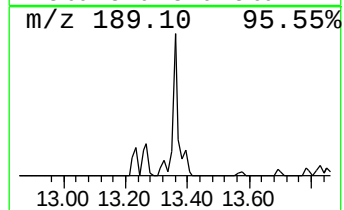
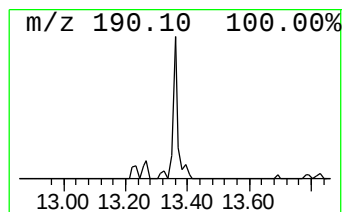
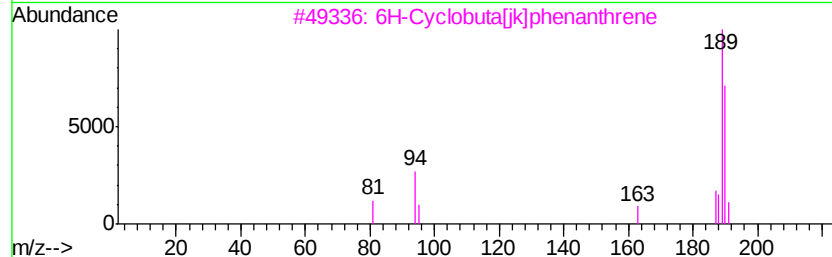
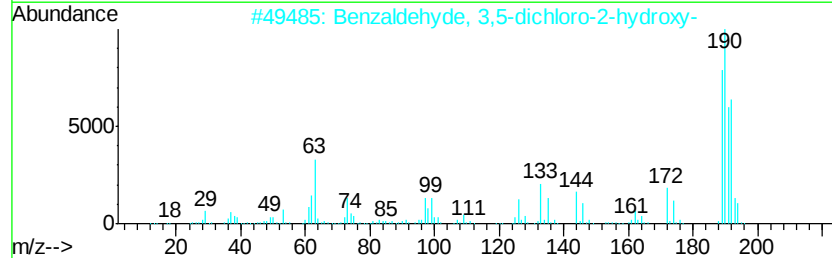
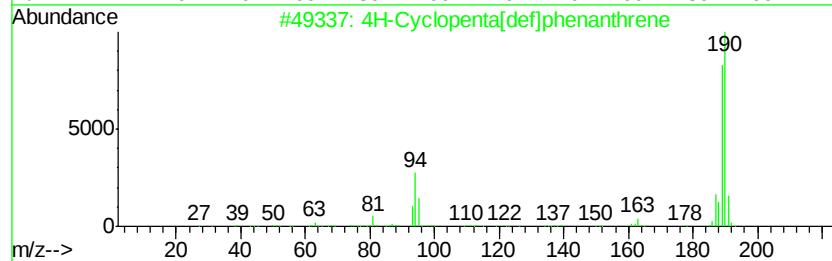
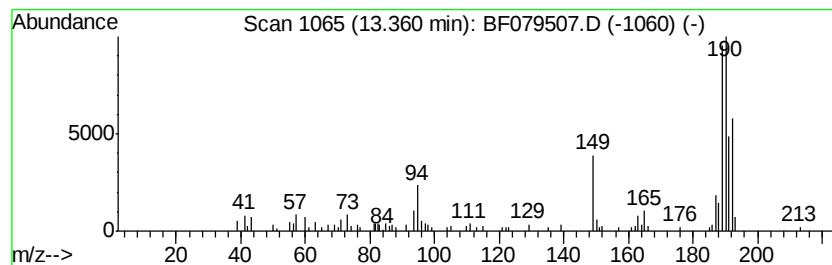
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 4H-Cyclopenta[def]phenanthrene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.36	5.53 ng	218301	Phenanthrene-d10	12.61

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	64
2		Benzaldehyde, 3,5-dichloro-2-hyd...	190	C7H4Cl2O2	000090-60-8	42
3		6H-Cyclobuta[ik]phenanthrene	190	C15H10	083469-43-6	38
4		Methyl diselenide	190	C2H6Se2	007101-31-7	38
5		Phenol, 4-(2-thienylmethyl)-	190	C11H10OS	091680-55-6	32



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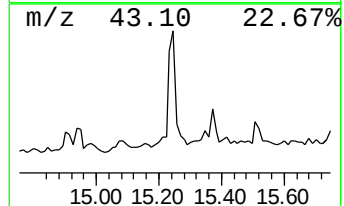
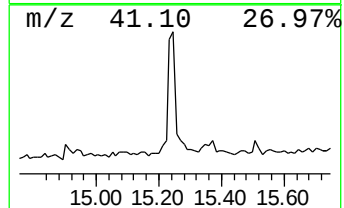
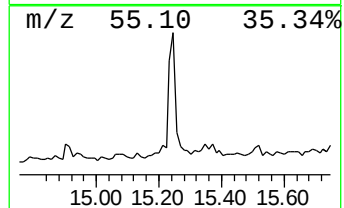
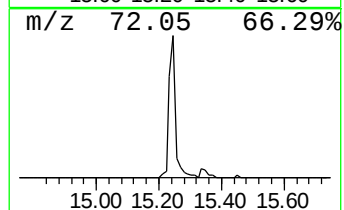
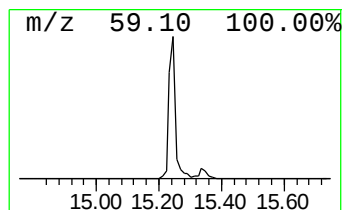
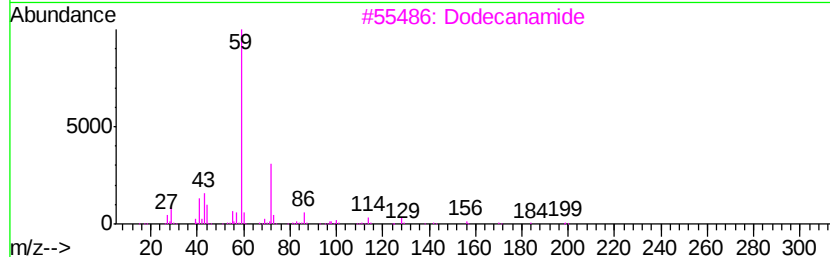
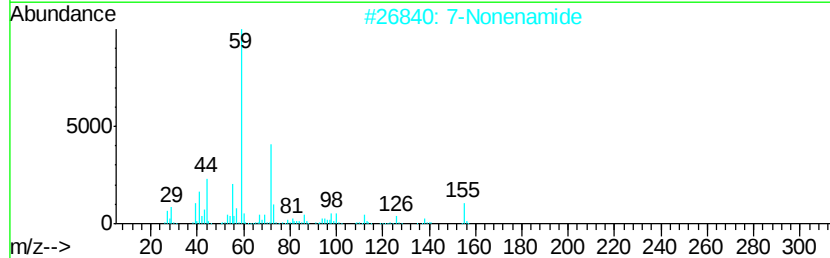
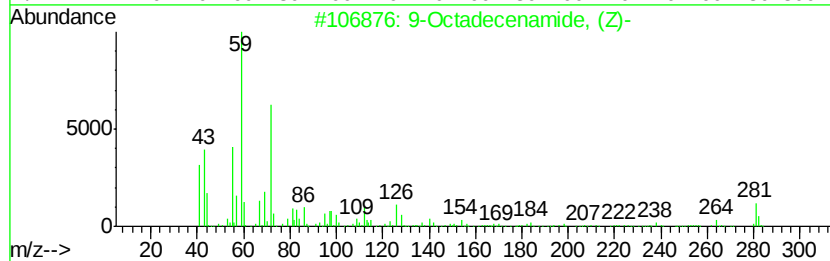
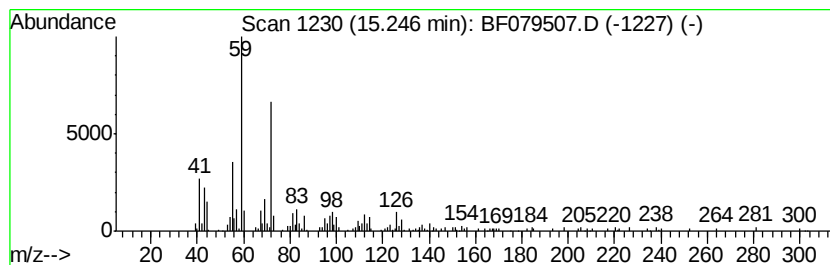
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 9-Octadecenamide, (Z)- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.25	3.41 ng	214358	Chrysene-d12	15.89

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	94
2		7-Nonenamide	155	C9H17NO	090949-53-4	78
3		Dodecanamide	199	C12H25NO	001120-16-7	53
4		Hexadecanamide	255	C16H33NO	000629-54-9	53
5		Tetradecanamide	227	C14H29NO	000638-58-4	53



Data Path : Z:\HPCHEM1\BNA F\DATA\BF052615\
Data File : BF079507.D
Acq On : 27 May 2015 7:08
Operator : TP/IZ
Sample : G2359-08
Misc :
ALS Vial : 30 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-30AS

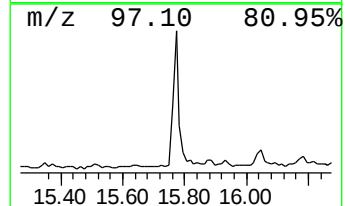
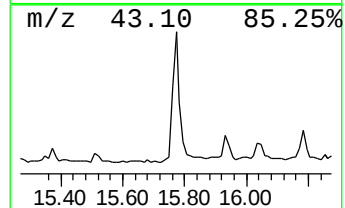
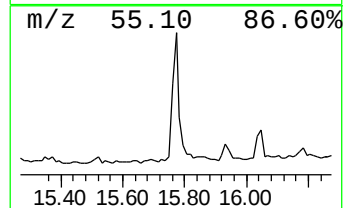
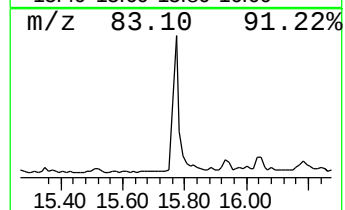
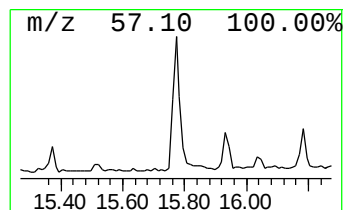
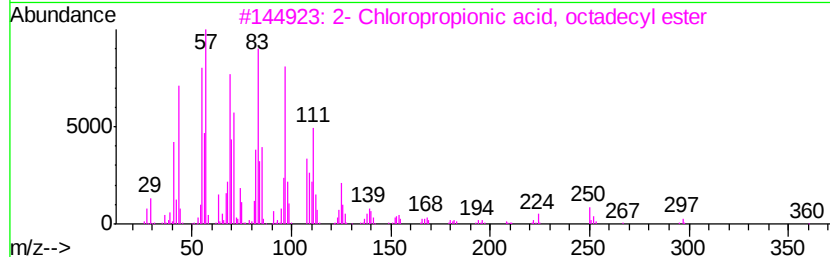
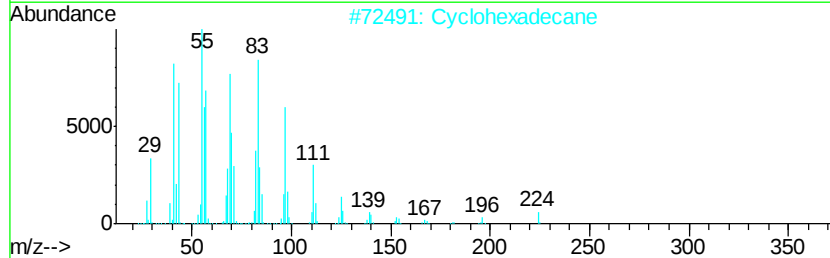
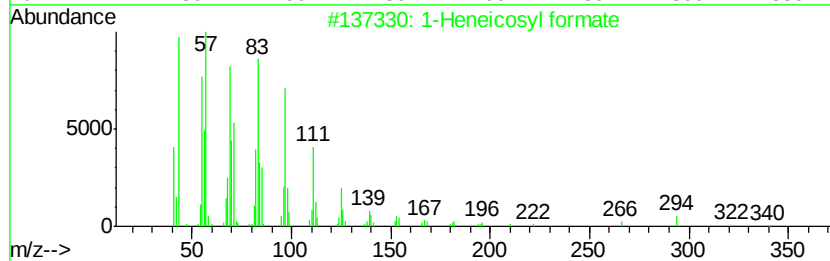
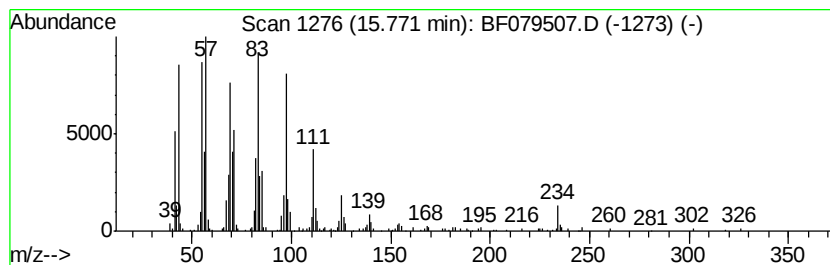
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 1-Heneicosyl formate Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.77	6.67 ng	419577	Chrysene-d12	15.89

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Heneicosyl formate	340	C22H44O2	077899-03-7	95
2		Cyclohexadecane	224	C16H32	000295-65-8	94
3		2- Chloropropionic acid, octadec...	360	C21H41ClO2	088104-31-8	94
4		Heptafluorobutanoic acid, heptad...	452	C21H35F7O2	1000282-97-3	94
5		Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	94



Data Path : Z:\HPCHEM1\BNA F\DATA\BF052615\
Data File : BF079507.D
Acq On : 27 May 2015 7:08
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Misc :
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BNA_F
ClientSampleId :
SB-30AS

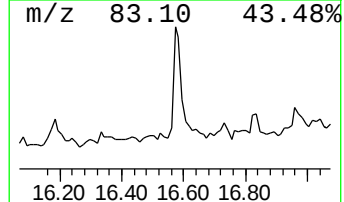
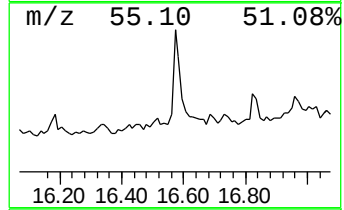
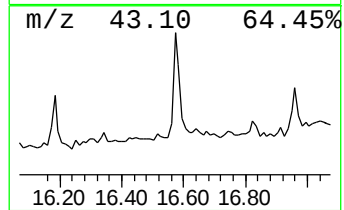
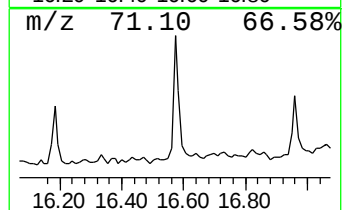
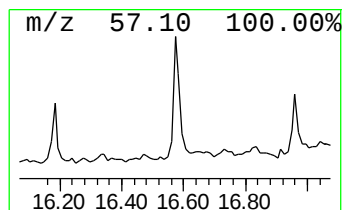
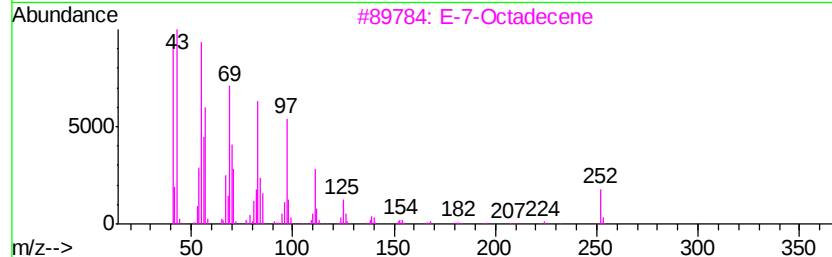
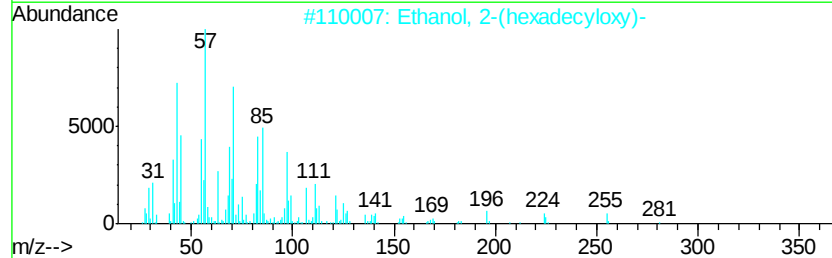
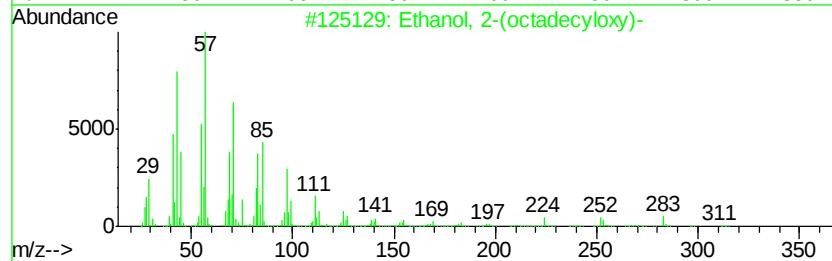
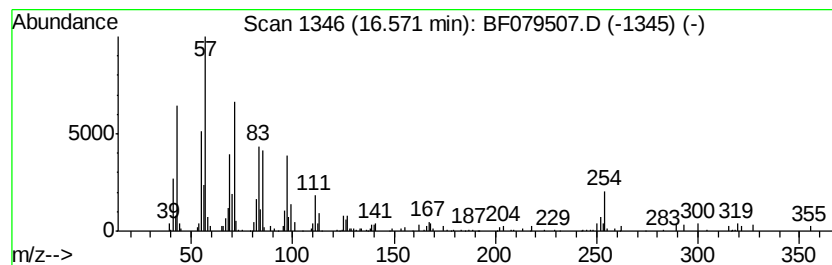
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 Ethanol, 2-(octadecyloxy)- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.57	2.35 ng	148074	Chrysene-d12	15.89

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Ethanol, 2-(octadecyloxy)-	314	C20H42O2	002136-72-3	70
2		Ethanol, 2-(hexadecyloxy)-	286	C18H38O2	002136-71-2	52
3		E-7-Octadecene	252	C18H36	1000130-92-0	46
4		Octadecane	254	C18H38	000593-45-3	45
5		Cyclotetracosane	336	C24H48	000297-03-0	43



Data Path : Z:\HPCHEM1\BNA F\DATA\BF052615\
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Acq On : 27 May 2015 7:08
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Misc :
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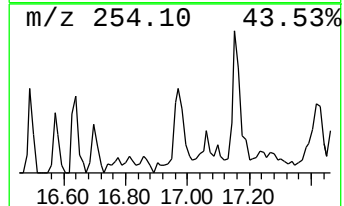
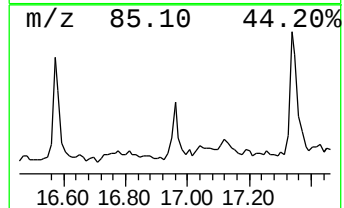
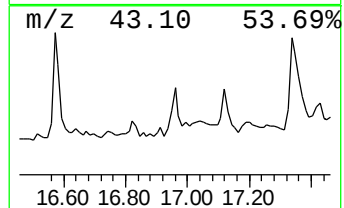
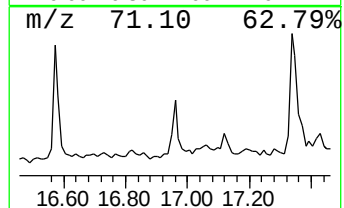
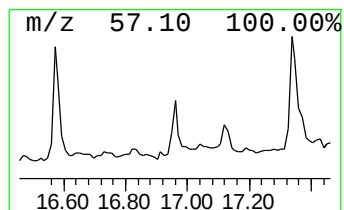
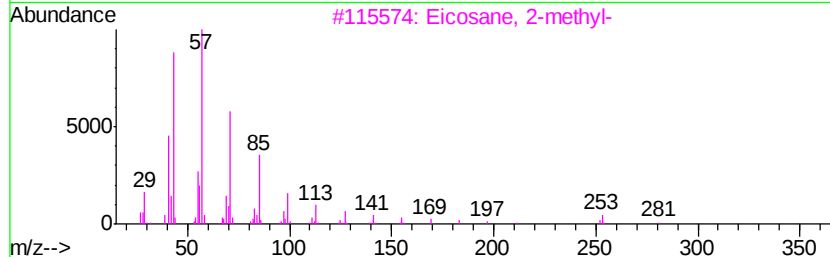
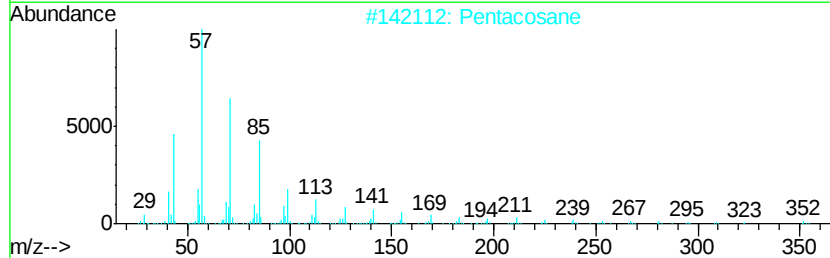
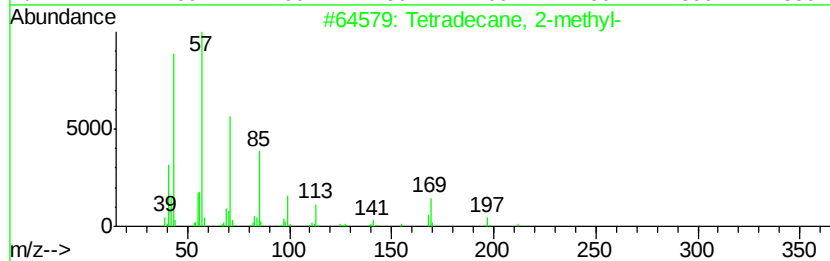
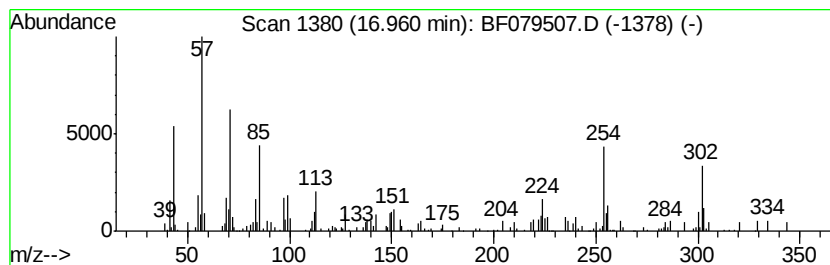
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 unknown16.96 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.96	3.45 ng	156882	Perylene-d12	17.59

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetradecane, 2-methyl-	212	C15H32	001560-95-8	35
2		Pentacosane	352	C25H52	000629-99-2	35
3		Eicosane, 2-methyl-	296	C21H44	001560-84-5	35
4		Eicosane	282	C20H42	000112-95-8	30
5		Nonadecane	268	C19H40	000629-92-5	27



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF052615\
Data File : BF079507.D
Acq On : 27 May 2015 7:08
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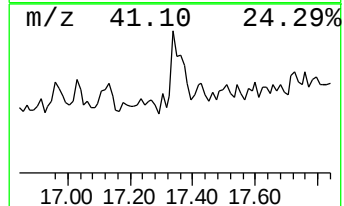
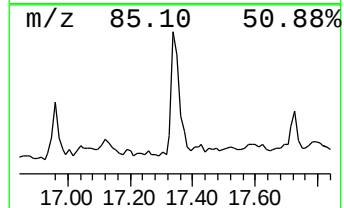
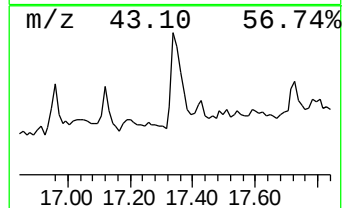
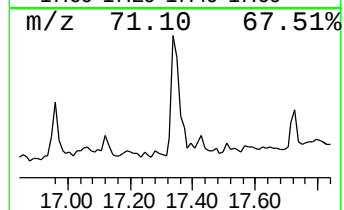
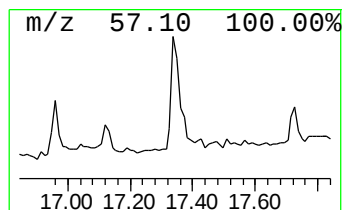
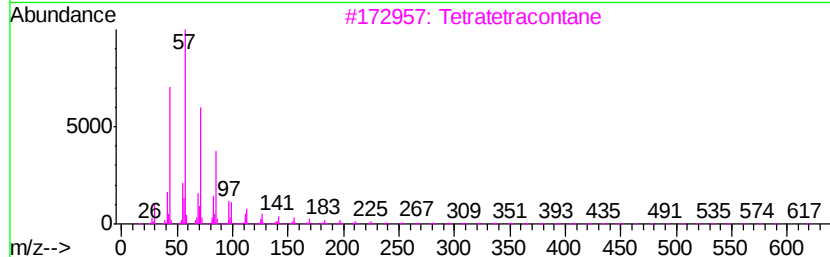
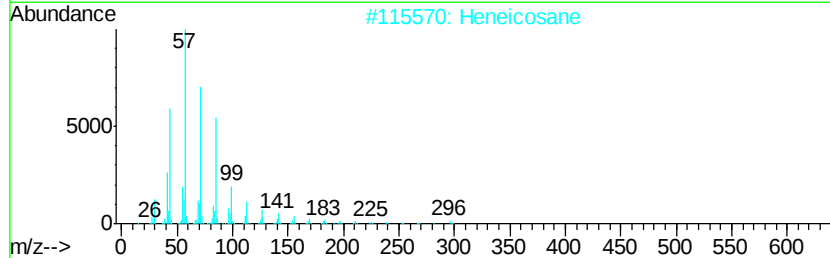
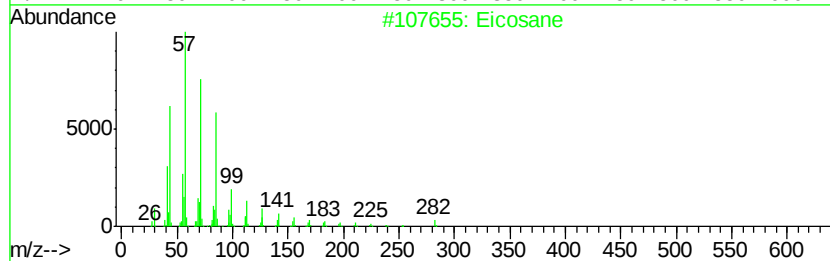
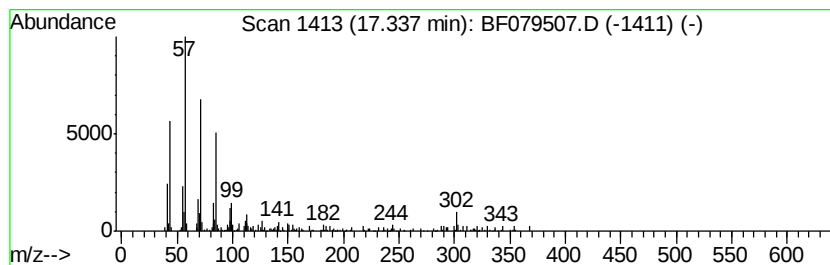
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 14 Eicosane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.34	5.17 ng	235345	Perylene-d12	17.59

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	94
2		Heneicosane	296	C21H44	000629-94-7	83
3		Tetratetracontane	619	C44H90	007098-22-8	83
4		Hexatriacontane	507	C36H74	000630-06-8	83
5		Tricosane, 2-methyl-	338	C24H50	001928-30-9	80



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF052615\
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Acq On : 27 May 2015 7:08
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Sample : G2359-08
Misc :
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Instrument :
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ClientSampleId :
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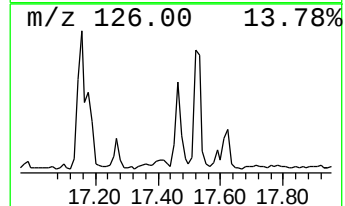
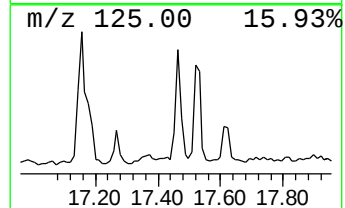
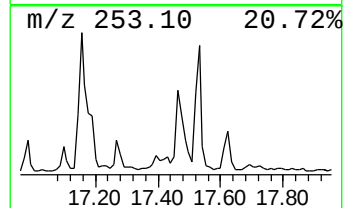
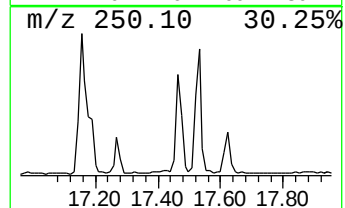
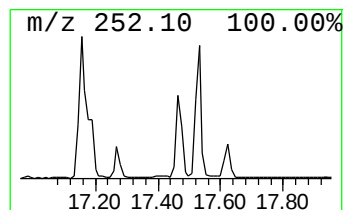
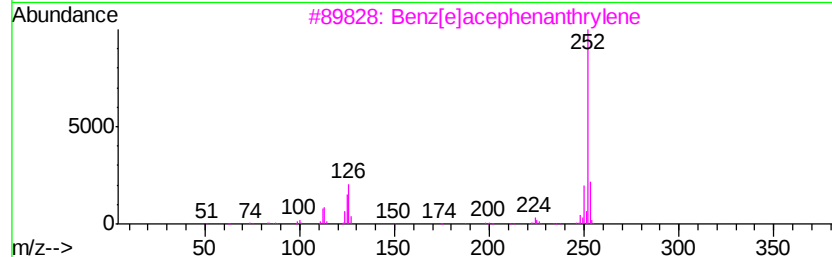
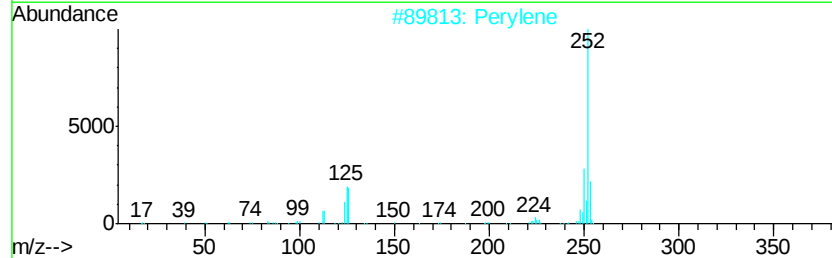
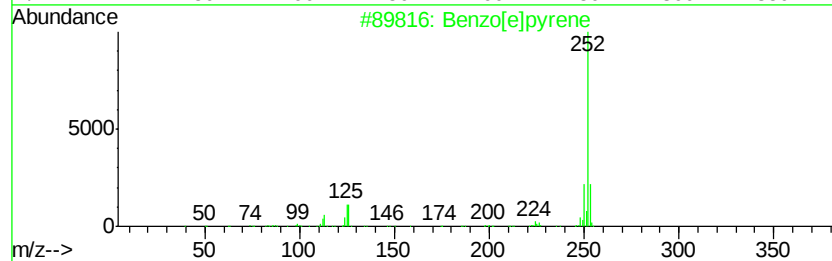
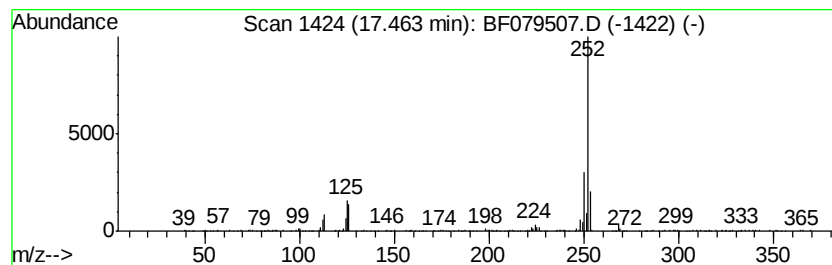
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 15 Benzo[e]pyrene Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.46	4.78 ng	217481	Perylene-d12	17.59

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[e]pyrene	252	C20H12	000192-97-2	94
2			Perylene	252	C20H12	000198-55-0	94
3			Benzo[e]acephenanthrylene	252	C20H12	000205-99-2	93
4			Benzo[k]fluoranthene	252	C20H12	000207-08-9	93
5			Benzo[a]pyrene	252	C20H12	000050-32-8	92



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF052615\
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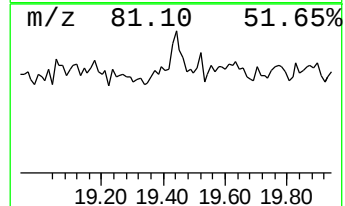
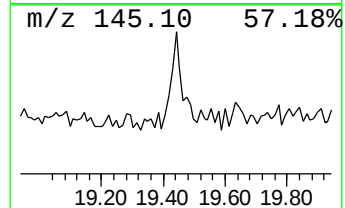
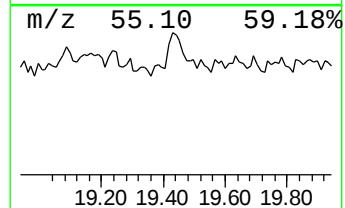
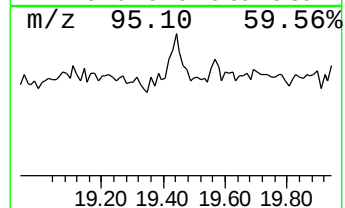
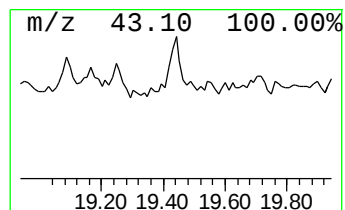
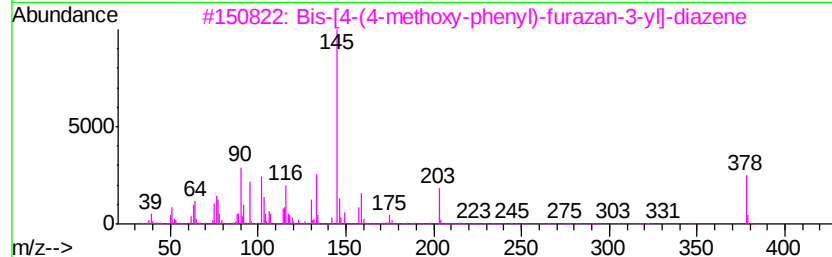
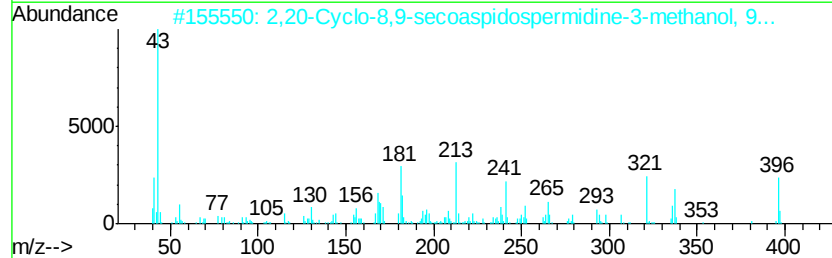
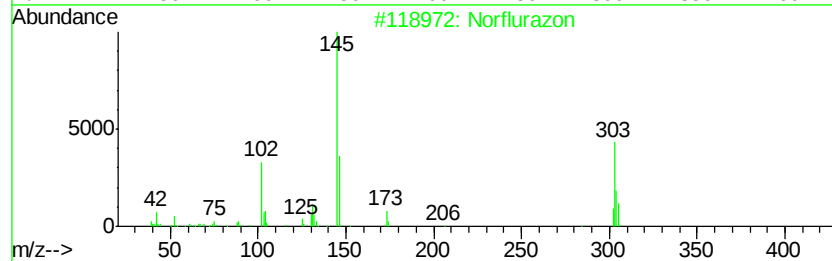
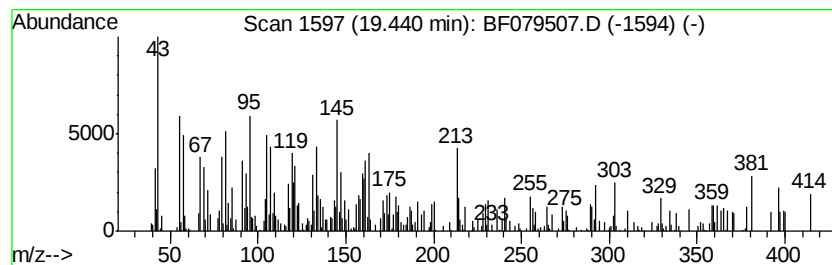
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 17 unknown19.44 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.44	4.00 ng	181817	Perylene-d12	17.59

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Norflurazon	303	C12H9ClF3N3O	027314-13-2	11
2		2,20-Cyclo-8,9-secoaspidospermid...	396	C24H32N2O3	056145-37-0	10
3		Bis-[4-(4-methoxy-phenyl)-furaza...	378	C18H14N6O4	1000296-30-1	9
4		5-Cholesten-3.beta.-acetoxv-24-t...	460	C29H48O2S	1000253-09-1	9
5		22-Ketocholesterol	400	C27H44O2	1000127-06-3	9



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF052615\
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF052615.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-methoxy...	1.43	10.6	ng	231896	1	7.02	437910	20.0
2-Pentanone, 4-hy...	4.74	7.4	ng	161380	1	7.02	437910	20.0
4H-Cyclopenta[def...	13.36	5.5	ng	218301	4	12.61	789772	20.0
9-Octadecenamide,...	15.25	3.4	ng	214358	5	15.89	1258600	20.0
1-Heneicosyl formate	15.77	6.7	ng	419577	5	15.89	1258600	20.0
Ethanol, 2-(octad...	16.57	2.4	ng	148074	5	15.89	1258600	20.0
unknown16.96	16.96	3.5	ng	156882	6	17.59	910085	20.0
Eicosane	17.34	5.2	ng	235345	6	17.59	910085	20.0
Benzo[e]pyrene	17.46	4.8	ng	217481	6	17.59	910085	20.0
unknown19.44	19.44	4.0	ng	181817	6	17.59	910085	20.0