

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG071915\
 Data File : BG017978.D
 Acq On : 20 Jul 2015 3:37
 Operator : TP/UM
 Sample : G3009-01
 Misc :
 ALS Vial : 54 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 MW-01-071515

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_G\METHODS\8270-BG071715.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	2.788	13	17	25	rVV	54334	83814	1.31%	0.248%
2	2.859	25	29	40	rVB	1238878	1593487	24.90%	4.718%
3	4.733	343	348	358	rVB	60101	88237	1.38%	0.261%
4	5.191	419	426	434	rBV	678276	964645	15.07%	2.856%
5	6.766	685	694	704	rBV	433937	675425	10.55%	2.000%
6	7.118	747	754	762	rBV	1280567	2031063	31.74%	6.013%
7	7.583	826	833	840	rBV	354636	544285	8.51%	1.611%
8	7.900	880	887	894	rBV	1161069	1868206	29.19%	5.531%
9	8.681	1015	1020	1024	rBV	61928	99639	1.56%	0.295%
10	8.740	1024	1030	1043	rVB	919398	1605611	25.09%	4.754%
11	9.204	1105	1109	1116	rVB	41732	64454	1.01%	0.191%
12	9.569	1163	1171	1178	rBV6	33976	75497	1.18%	0.224%
13	9.768	1200	1205	1214	rVB3	105040	196916	3.08%	0.583%
14	10.362	1295	1306	1312	rBV	519788	905512	14.15%	2.681%
15	10.420	1312	1316	1322	rVV5	38546	77426	1.21%	0.229%
16	10.479	1322	1326	1333	rVB4	39542	70238	1.10%	0.208%
17	10.832	1374	1386	1391	rVB	53559	106172	1.66%	0.314%
18	10.943	1398	1405	1411	rBV2	58759	101271	1.58%	0.300%
19	11.560	1503	1510	1515	rBV2	47061	78853	1.23%	0.233%
20	11.925	1566	1572	1578	rVB2	80399	136823	2.14%	0.405%
21	12.259	1624	1629	1633	rBV	57637	89654	1.40%	0.265%
22	12.859	1724	1731	1737	rBV	2641773	3982762	62.24%	11.792%
23	13.705	1871	1875	1879	rBV	55276	79297	1.24%	0.235%
24	14.245	1957	1967	1974	rBV2	795523	1312181	20.51%	3.885%
25	14.704	2038	2045	2055	rBV	435410	739537	11.56%	2.190%
26	14.851	2065	2070	2077	rBV	88587	178719	2.79%	0.529%
27	15.614	2195	2200	2209	rBV3	110912	238173	3.72%	0.705%
28	15.738	2215	2221	2236	rVB	2281777	3409815	53.29%	10.096%
29	16.995	2429	2435	2441	rBV2	1090484	1555755	24.31%	4.606%
30	17.912	2586	2591	2599	rBV	163174	260160	4.07%	0.770%
31	19.204	2806	2811	2819	rBV5	65267	119934	1.87%	0.355%
32	19.457	2851	2854	2863	rVB	56344	88039	1.38%	0.261%
33	19.645	2880	2886	2891	rBV	5044303	6399115	100.00%	18.946%
34	20.920	3100	3103	3111	rVB	159825	203842	3.19%	0.604%

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ClientSampleId :
MW-01-071515

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

35	21.196	3145	3150	3163	rVB	1391295	1776979	27.77%	5.261%
36	22.330	3338	3343	3349	rBV	113103	176773	2.76%	0.523%
37	23.411	3520	3527	3540	rVB2	929578	1797201	28.09%	5.321%

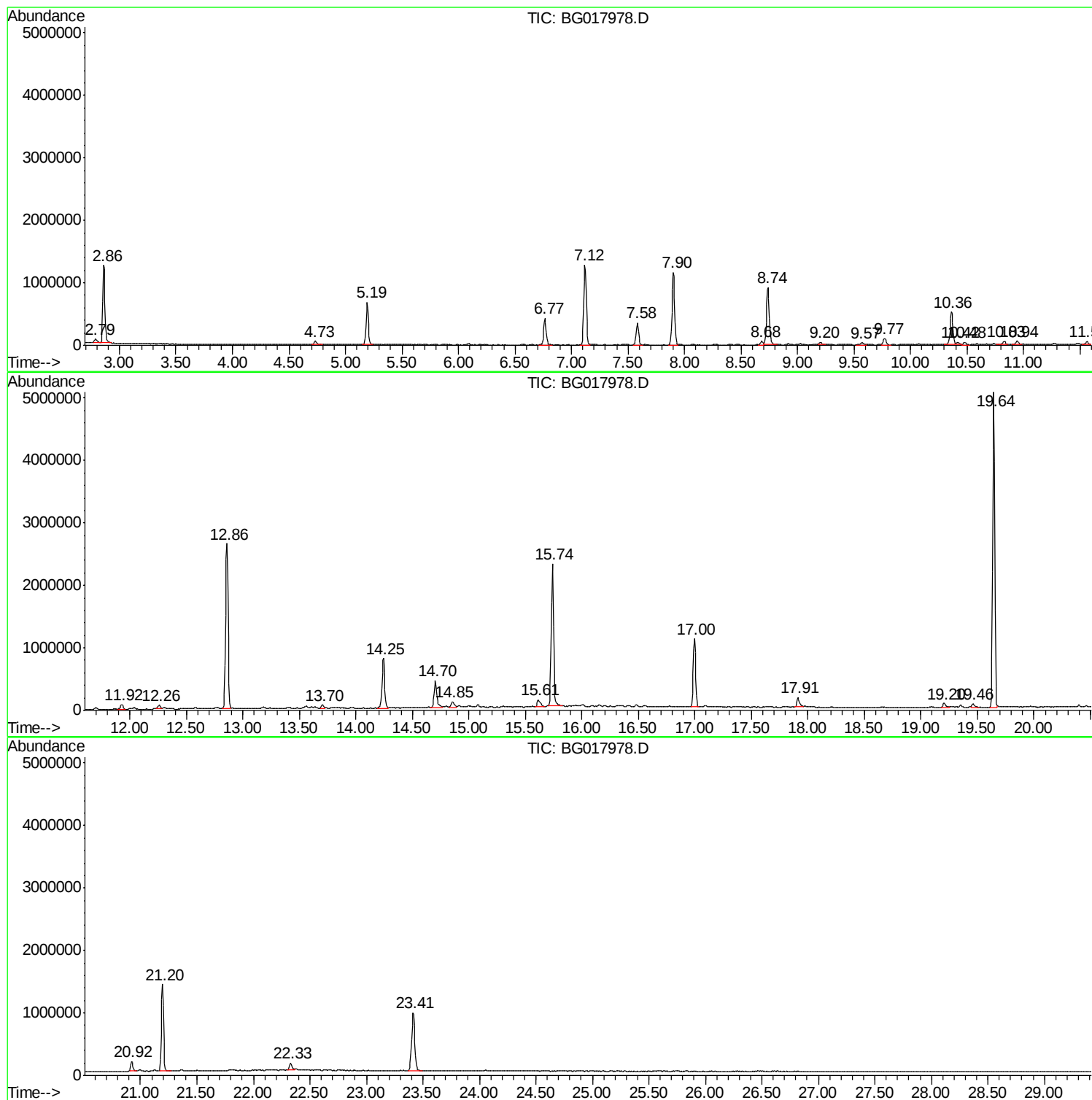
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Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG071715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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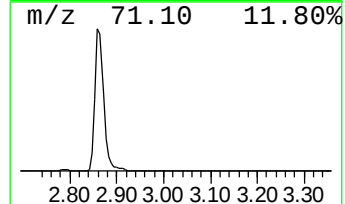
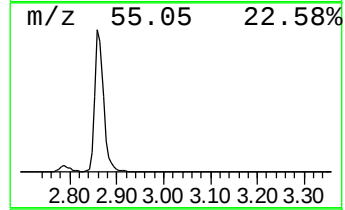
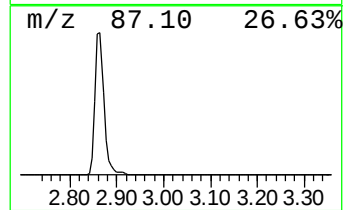
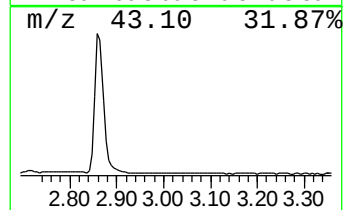
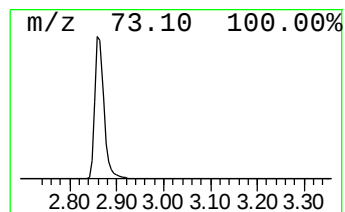
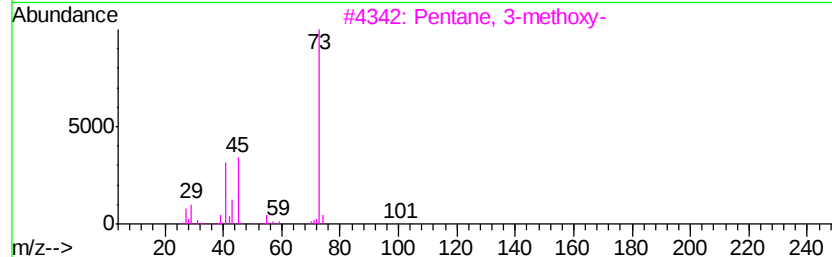
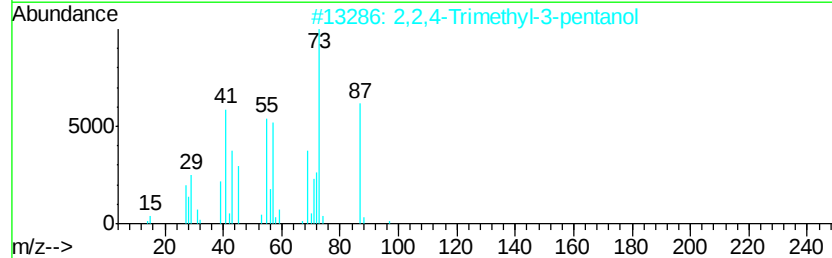
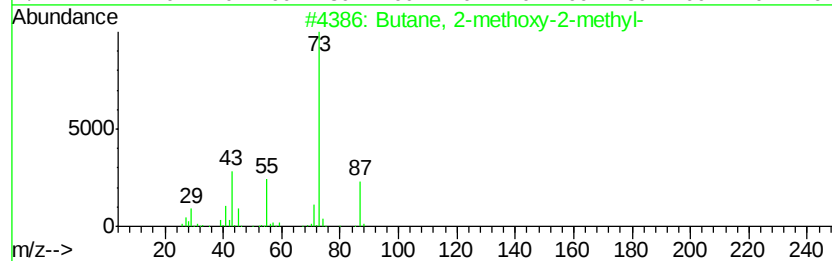
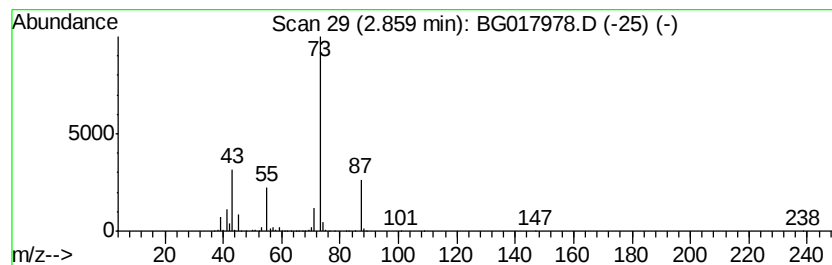
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG071715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 Butane, 2-methoxy-2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.86	58.55 ng	1593490	1,4-Dichlorobenzene-d4	7.58

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	86
2		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	64
3		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	25
4		1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	9
5		Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	9



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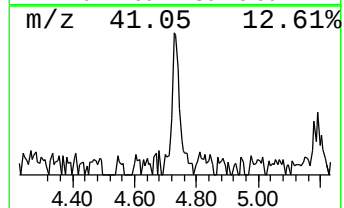
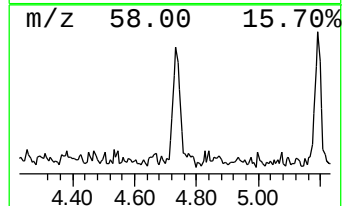
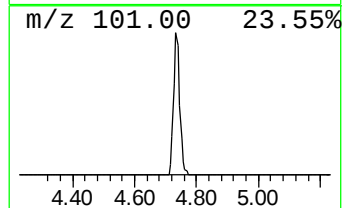
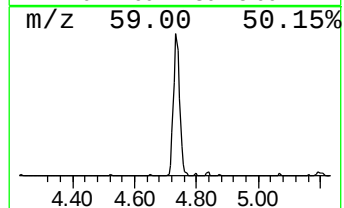
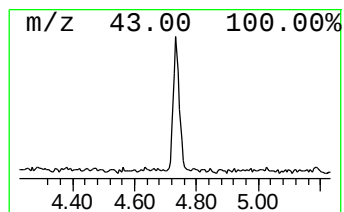
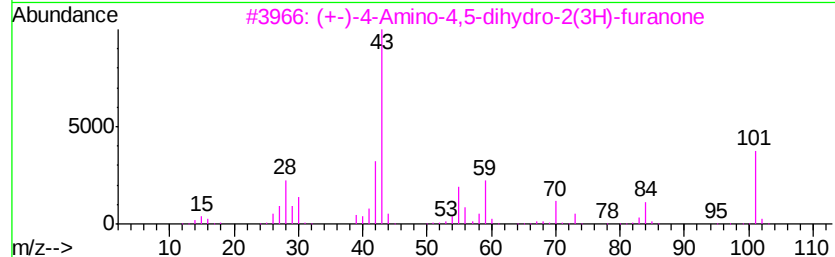
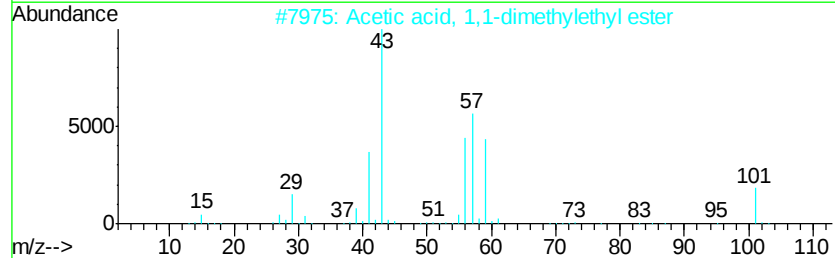
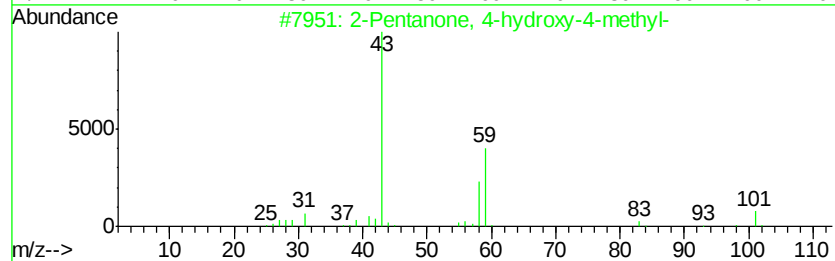
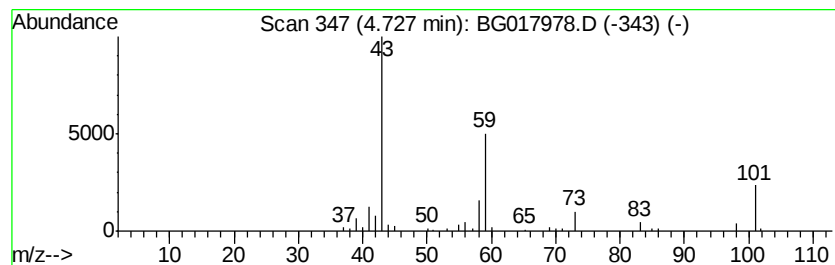
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.73	3.24 ng	88237	1,4-Dichlorobenzene-d4	7.58

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	45
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	38
3			(+)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	25
4			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	25
5			4-Ethyl-3-octanol	158	C10H22O	019781-26-1	23



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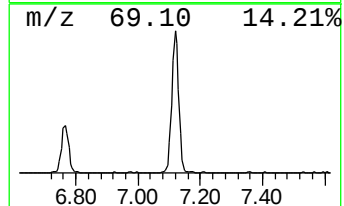
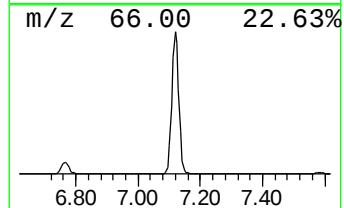
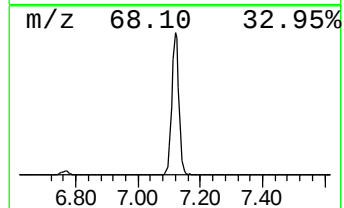
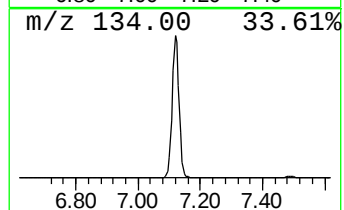
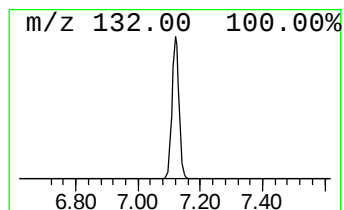
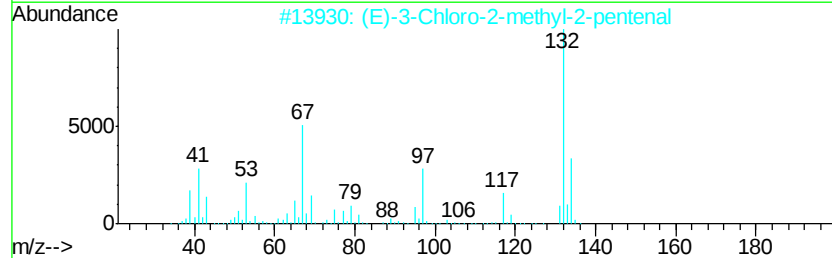
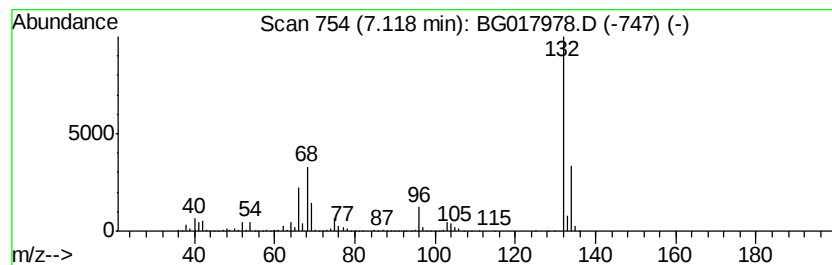
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 4 unknown7.12 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.12	74.63 ng	2031060	1,4-Dichlorobenzene-d4	7.58

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	35
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	25
3		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
4		5-Aminoindole	132	C8H8N2	005192-03-0	10
5		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9



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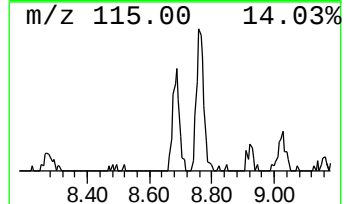
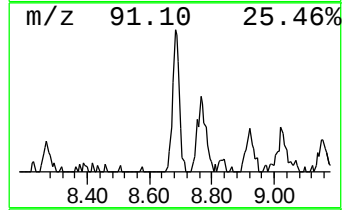
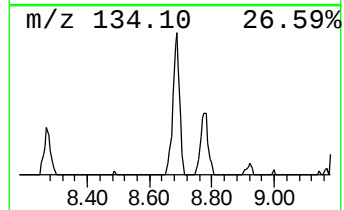
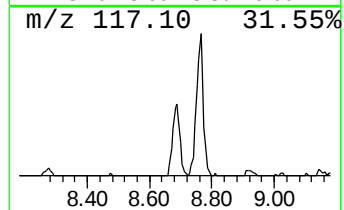
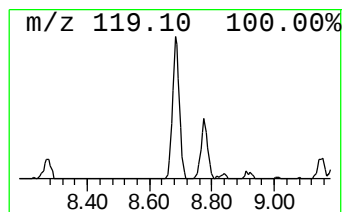
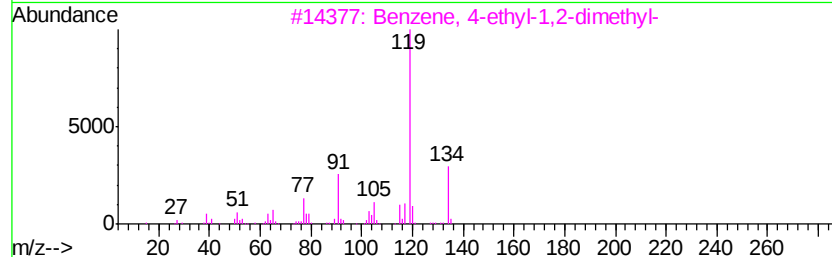
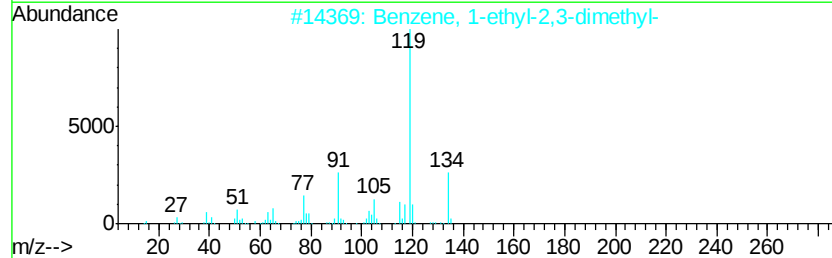
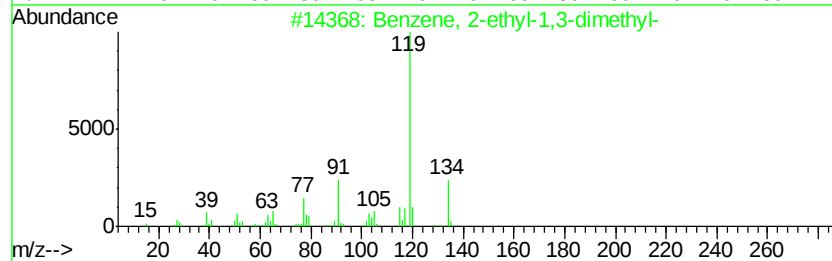
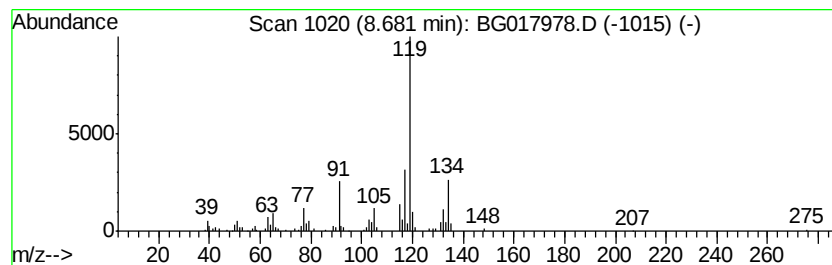
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Benzene, 2-ethyl-1,3-dimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.68	3.66 ng	99639	1,4-Dichlorobenzene-d4	7.58

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	93
2		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	93
3		Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	93
4		Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	93
5		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	81



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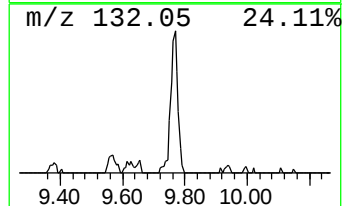
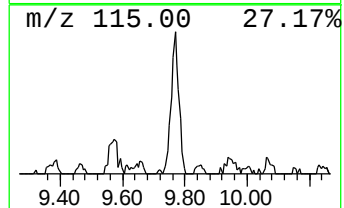
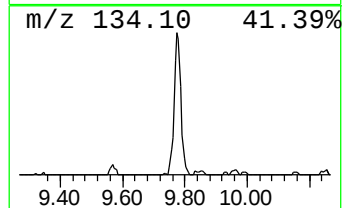
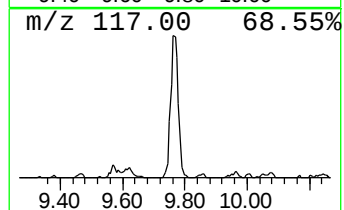
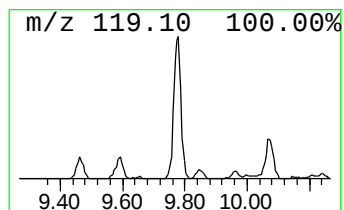
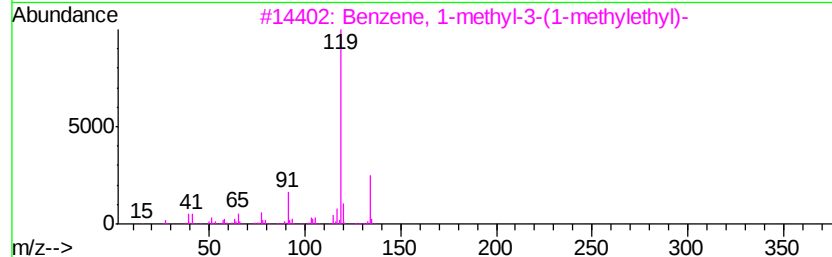
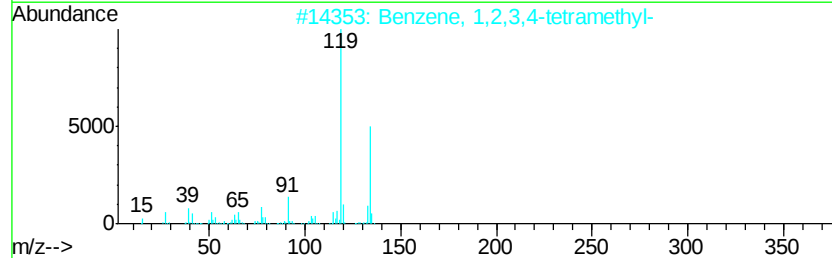
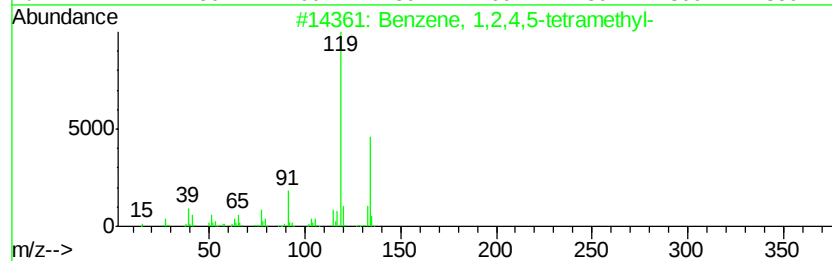
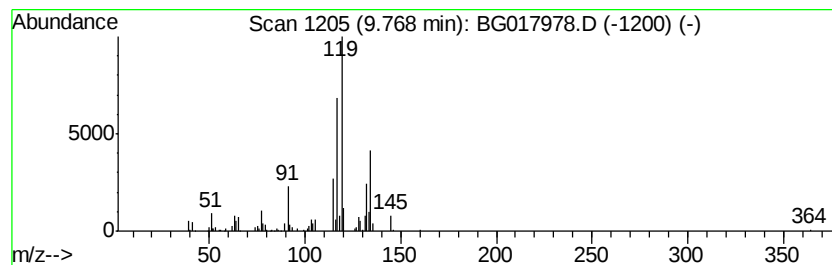
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.77	4.35 ng	196916	Naphthalene-d8	10.36

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	64
2		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	64
3		Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	60
4		Benzene, 1-methyl-4-(1-methyleth...	134	C10H14	000099-87-6	60
5		Benzene, 1-methyl-2-(1-methyleth...	134	C10H14	000527-84-4	60



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Sample : G3009-01
Misc :
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Instrument :
BNA_G
ClientSampleId :
MW-01-071515

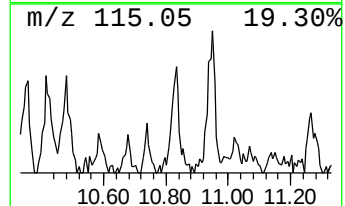
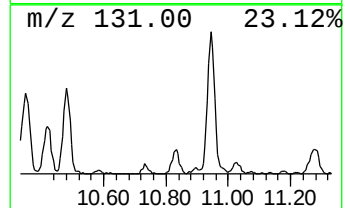
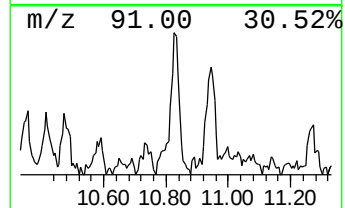
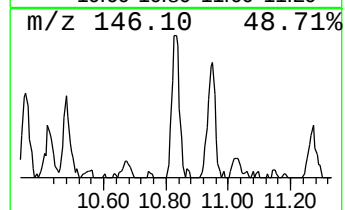
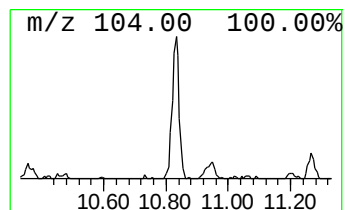
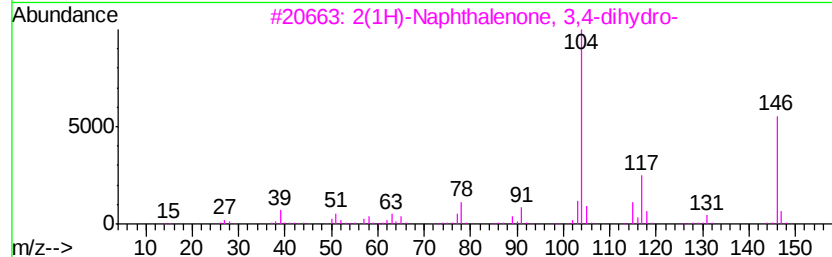
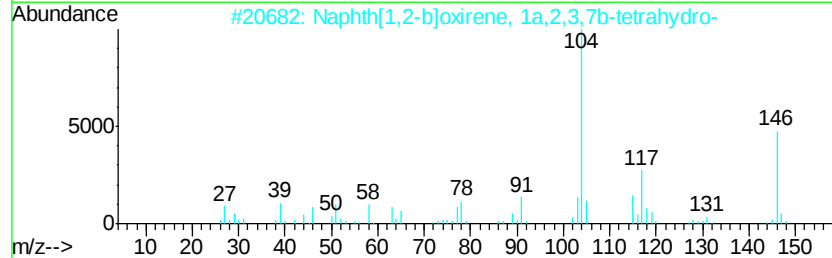
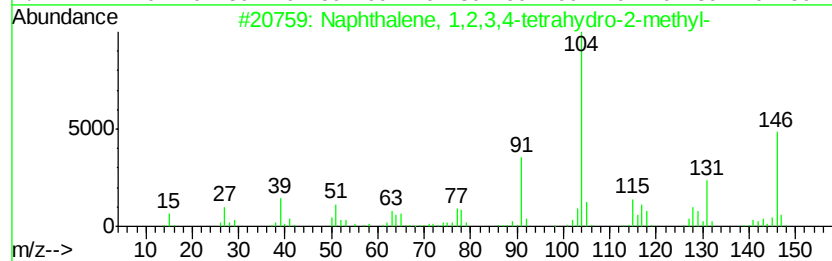
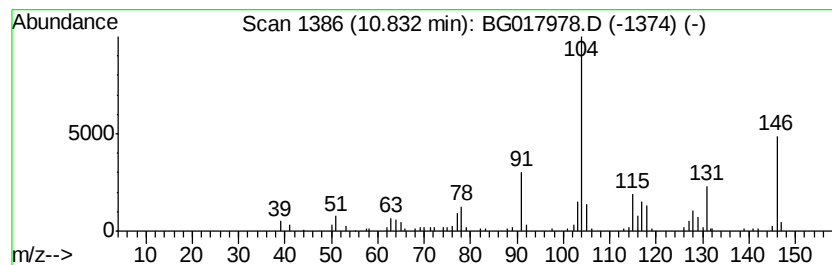
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Naphthalene, 1,2,3,4-tetra... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.83	2.35 ng	106172	Naphthalene-d8	10.36

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	003877-19-8	94
2		Naphth[1,2-b]oxirene, 1a,2,3,7b-...	146	C10H10O	002461-34-9	62
3		2(1H)-Naphthalenone, 3,4-dihydro-	146	C10H10O	000530-93-8	43
4		2-Methylbenzyl cyanide	131	C9H9N	022364-68-7	38
5		1,3-Propanediamine, 2-phenyl-	150	C9H14N2	055165-09-8	35



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG071915\
Data File : BG017978.D
Acq On : 20 Jul 2015 3:37
Operator : TP/UM
Sample : G3009-01
Misc :
ALS Vial : 54 Sample Multiplier: 1

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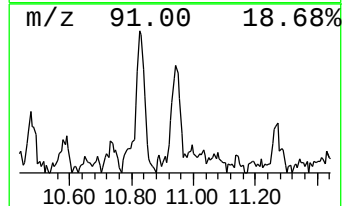
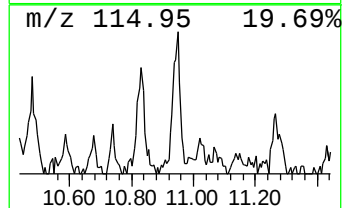
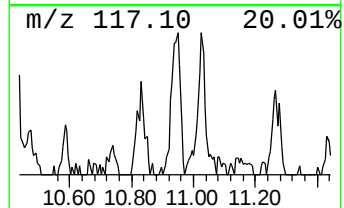
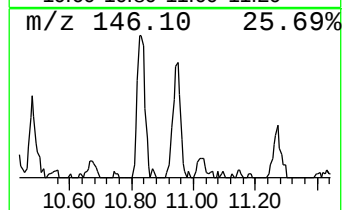
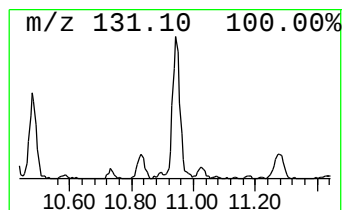
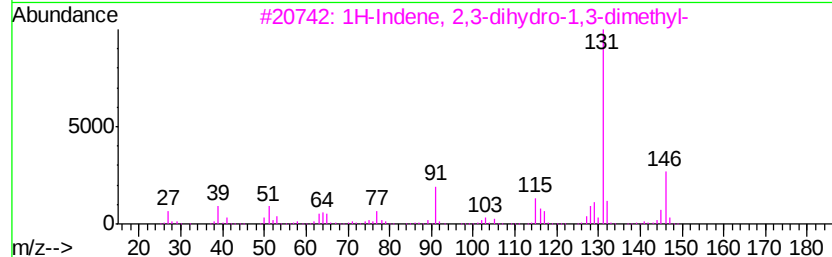
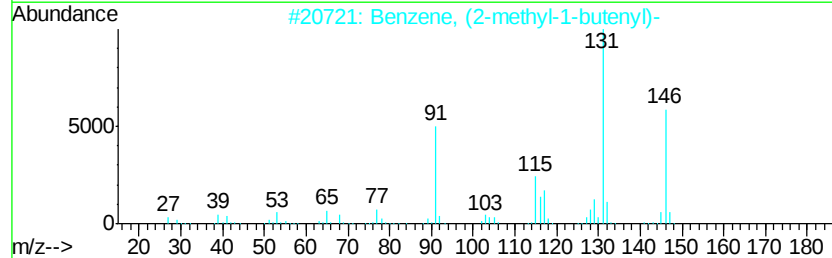
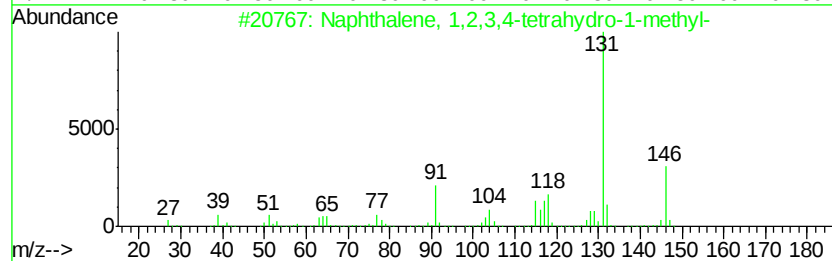
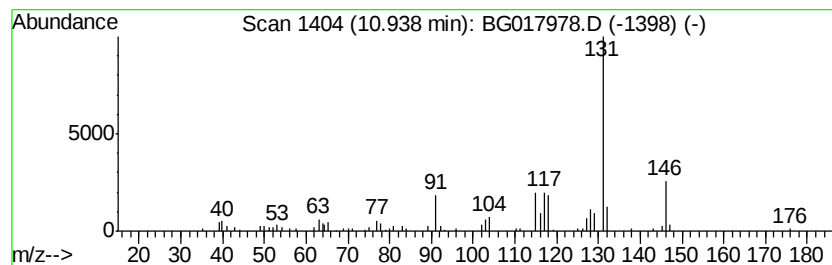
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Naphthalene, 1,2,3,4-tetra... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.94	2.24 ng	101271	Naphthalene-d8	10.36

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001559-81-5	94
2		Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	93
3		1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	90
4		1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	90
5		1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	87



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG071915\
Data File : BG017978.D
Acq On : 20 Jul 2015 3:37
Operator : TP/UM
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Misc :
ALS Vial : 54 Sample Multiplier: 1

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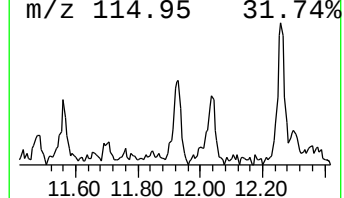
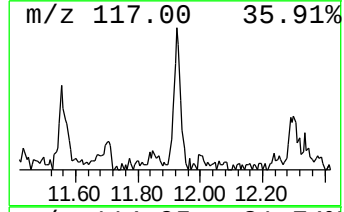
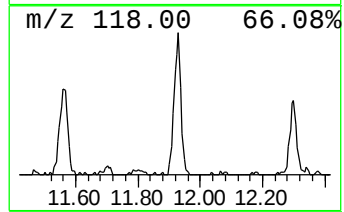
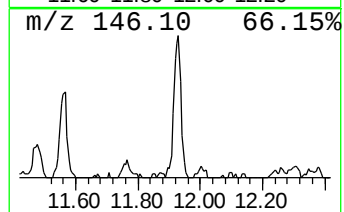
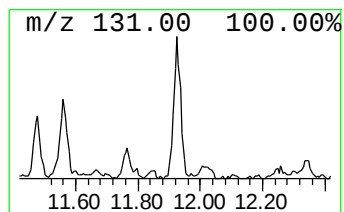
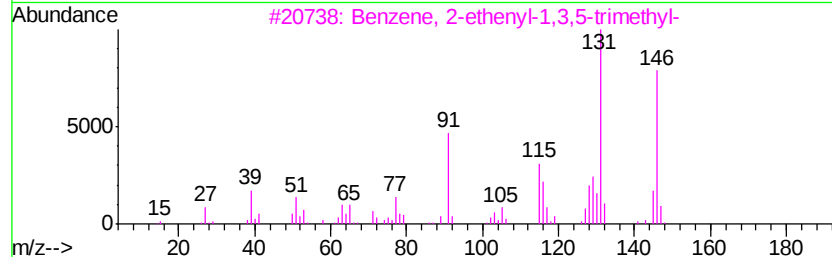
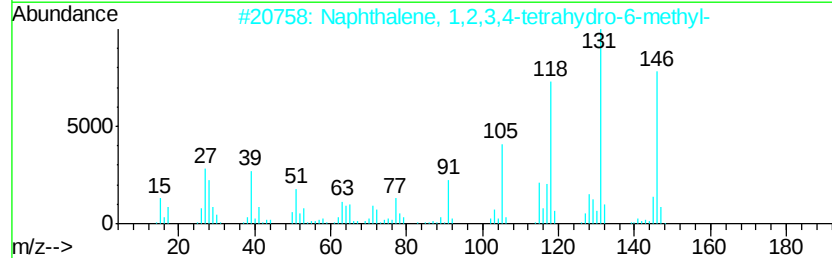
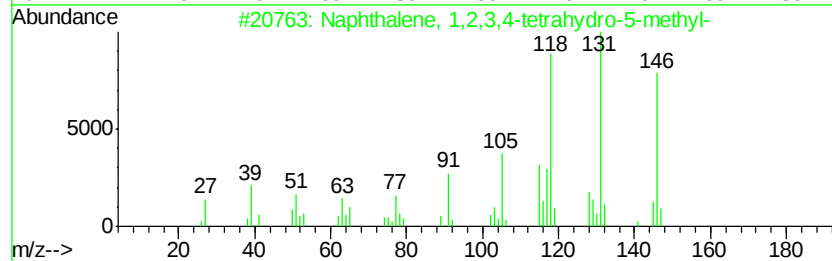
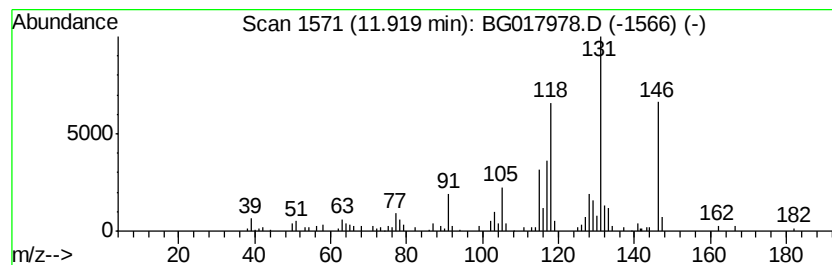
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Naphthalene, 1,2,3,4-tetra... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.92	3.02 ng	136823	Naphthalene-d8	10.36

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	95
2		Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	87
3		Benzene, 2-ethenyl-1,3,5-trimethyl-	146	C11H14	000769-25-5	86
4		Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	76
5		1H-Indene,2,3-dihydro-2,2-dimethyl-	146	C11H14	020836-11-7	70



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG071915\
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Operator : TP/UM
Sample : G3009-01
Misc :
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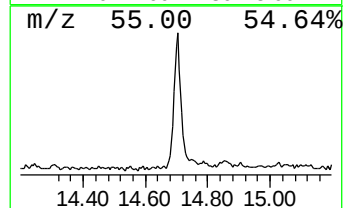
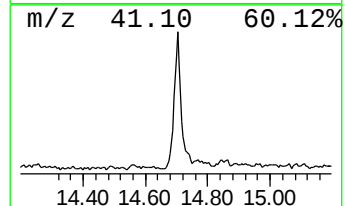
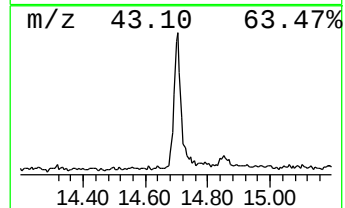
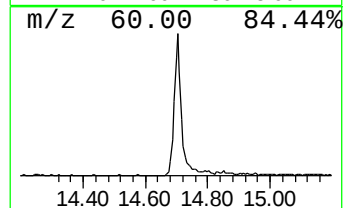
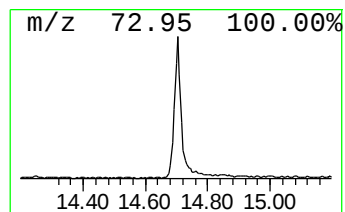
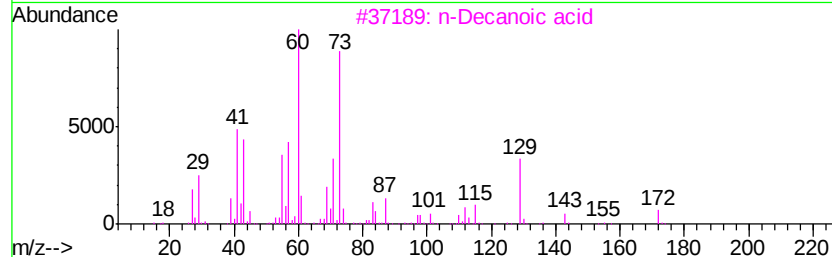
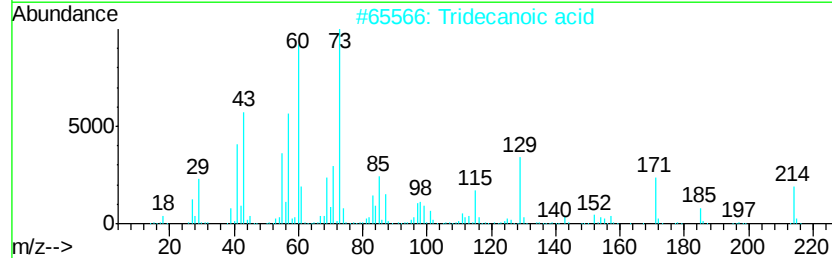
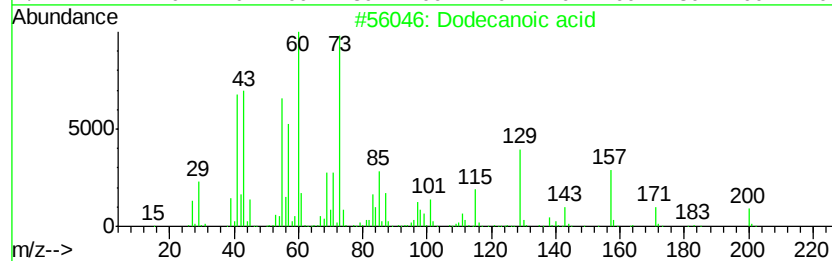
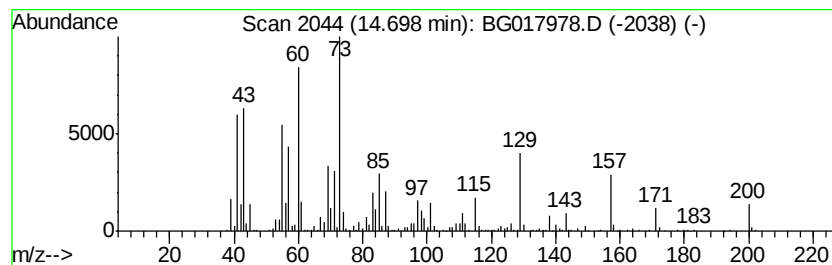
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Dodecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.70	11.27 ng	739537	Acenaphthene-d10	14.24
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Dodecanoic acid	200 C12H24O2	000143-07-7 99
2		Tridecanoic acid	214 C13H26O2	000638-53-9 59
3		n-Decanoic acid	172 C10H20O2	000334-48-5 53
4		Octanoic Acid	144 C8H16O2	000124-07-2 43
5		Nonanoic acid	158 C9H18O2	000112-05-0 43



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG071915\
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Acq On : 20 Jul 2015 3:37
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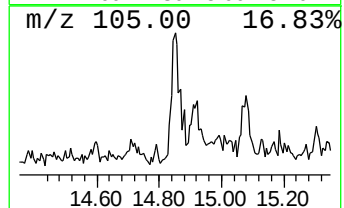
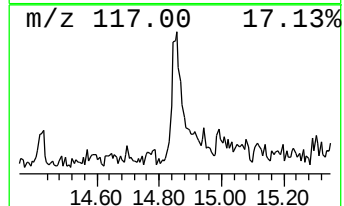
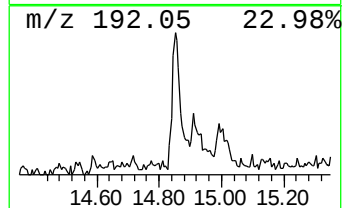
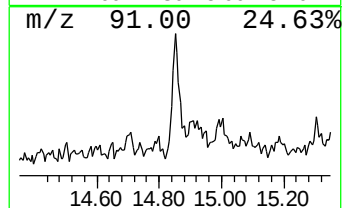
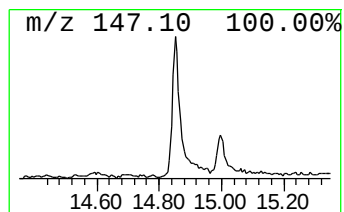
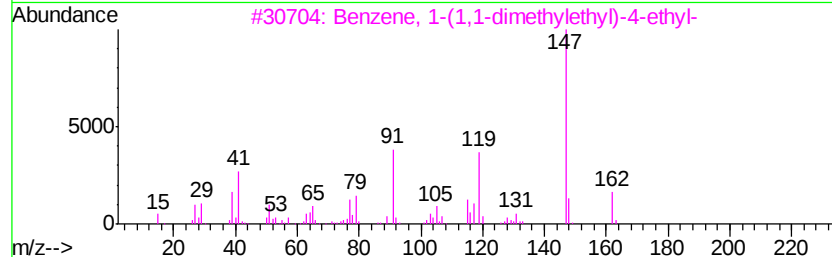
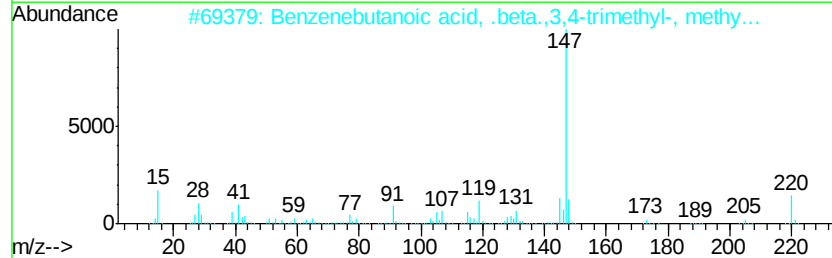
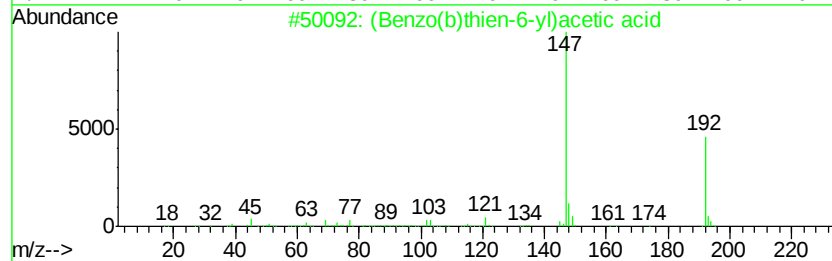
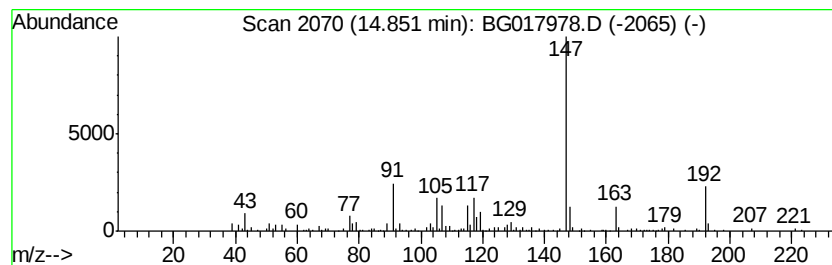
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 (Benzo(b)thien-6-yl)acetic ... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.85	2.72 ng	178719	Acenaphthene-d10	14.24
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		(Benzo(b)thien-6-yl)acetic acid	192 C10H8O2S	006177-87-3 53
2		Benzenebutanoic acid, .beta.,3,4...	220 C14H20O2	069855-51-2 50
3		Benzene, 1-(1,1-dimethylethyl)-4...	162 C12H18	007364-19-4 47
4		Benzoic acid, 4-propyl-, 4-cyano...	265 C17H15NO2	056131-49-8 47
5		2,4,6-Trifluoroaniline	147 C6H4F3N	000363-81-5 47



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG071915\
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Misc :
ALS Vial : 54 Sample Multiplier: 1

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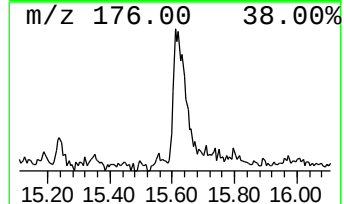
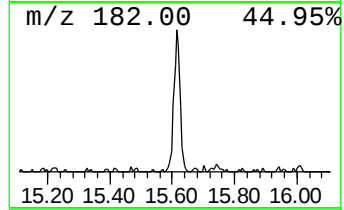
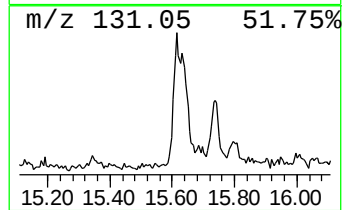
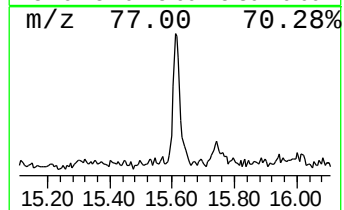
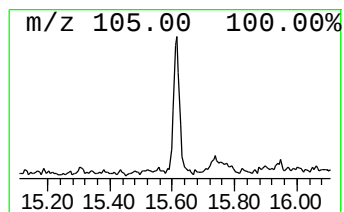
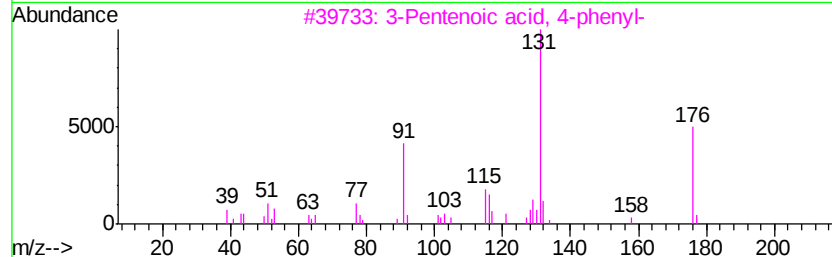
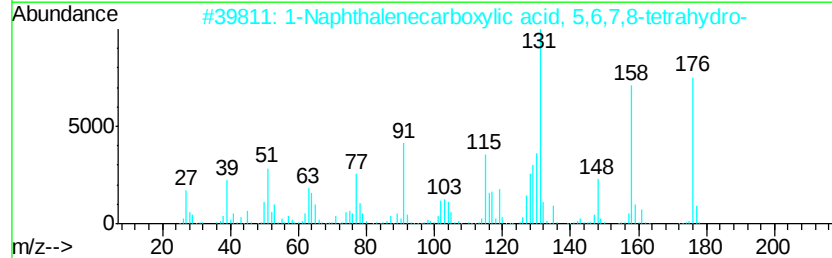
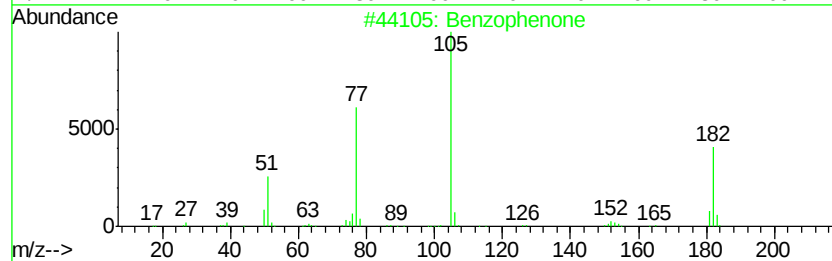
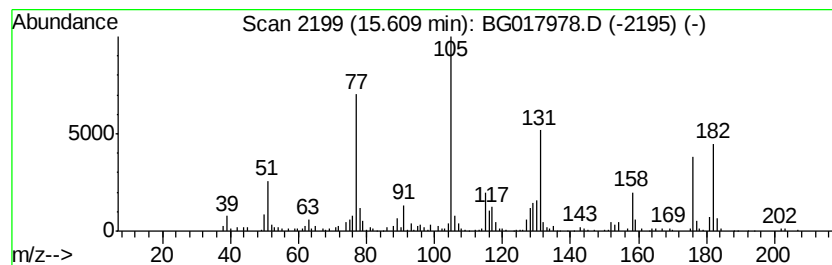
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Benzophenone Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.61	3.63 ng	238173	Acenaphthene-d10	14.24

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzophenone	182	C13H10O	000119-61-9	52
2		1-Naphthalenecarboxylic acid, 5,...	176	C11H12O2	004242-18-6	38
3		3-Pentenoic acid, 4-phenyl-	176	C11H12O2	053774-19-9	30
4		Benzeneacetonitrile, .alpha.-oxo-	131	C8H5NO	000613-90-1	27
5		2-Propenoic acid, 3-phenyl-, eth...	176	C11H12O2	000103-36-6	18



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG071915\
Data File : BG017978.D
Acq On : 20 Jul 2015 3:37
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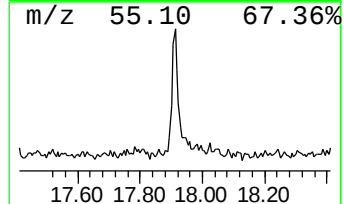
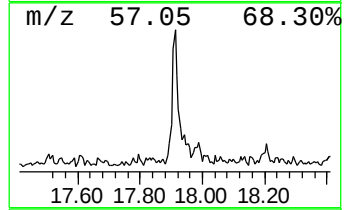
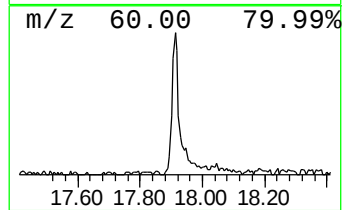
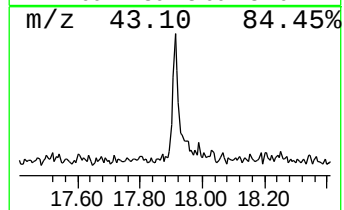
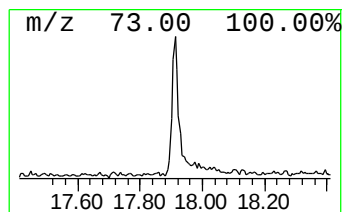
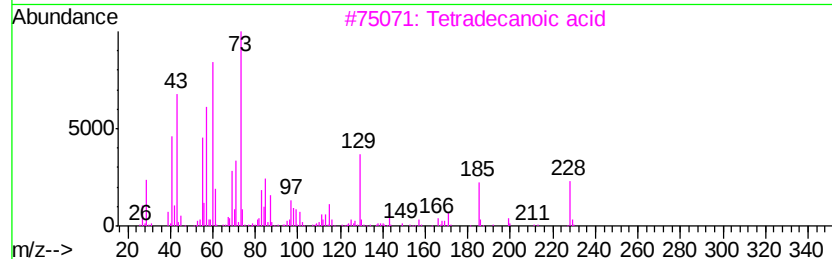
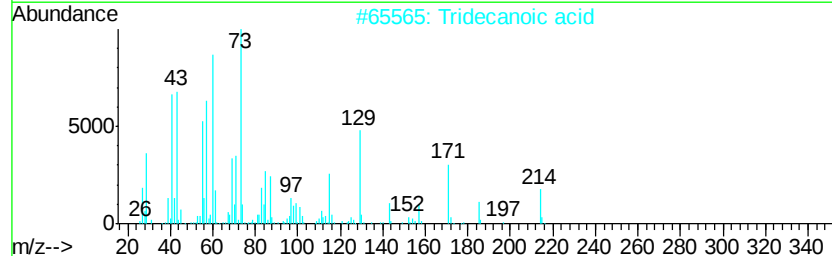
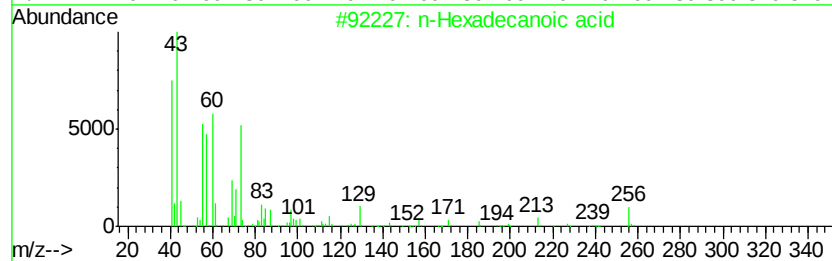
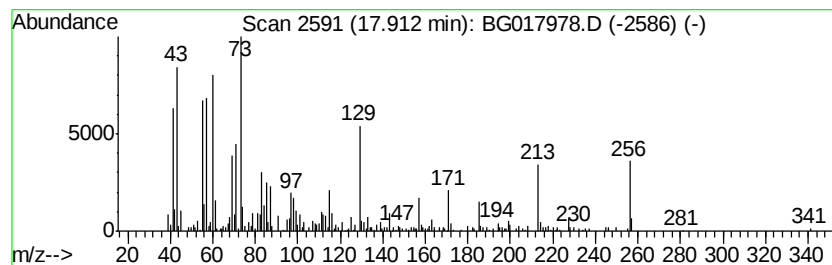
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG071715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 n-Hexadecanoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.91	3.34 ng	260160	Phenanthrene-d10	17.00

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95
2		Tridecanoic acid	214	C13H26O2	000638-53-9	94
3		Tetradecanoic acid	228	C14H28O2	000544-63-8	94
4		Pentadecanoic acid	242	C15H30O2	001002-84-2	76
5		n-Decanoic acid	172	C10H20O2	000334-48-5	64



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG071915\
Data File : BG017978.D
Acq On : 20 Jul 2015 3:37
Operator : TP/UM
Sample : G3009-01
Misc :
ALS Vial : 54 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
MW-01-071515

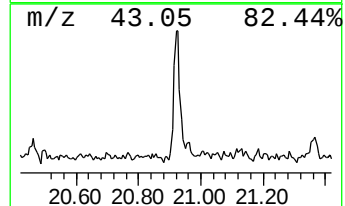
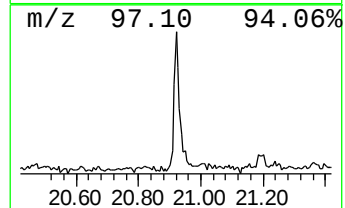
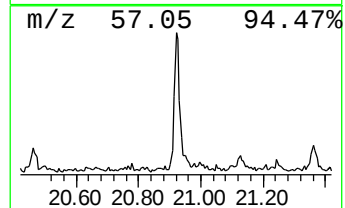
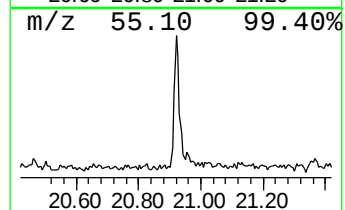
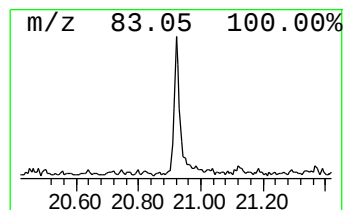
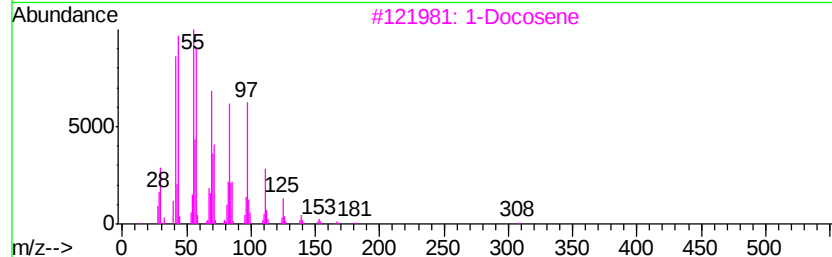
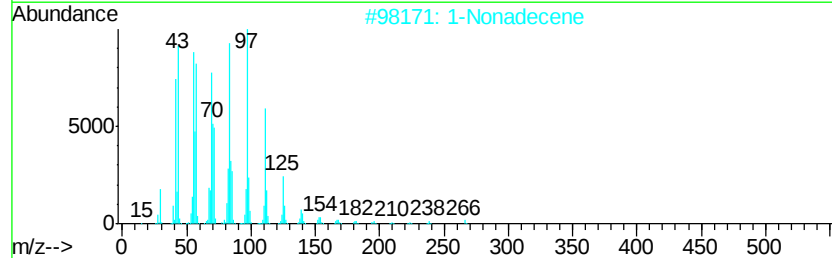
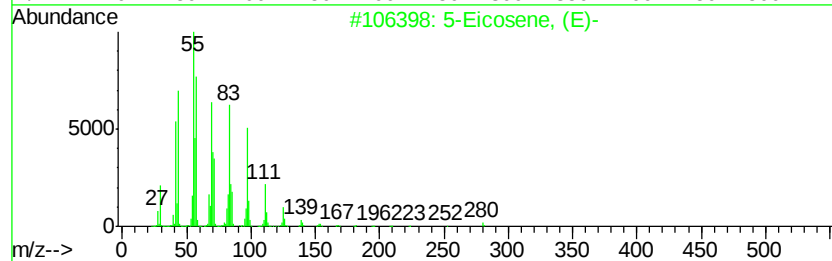
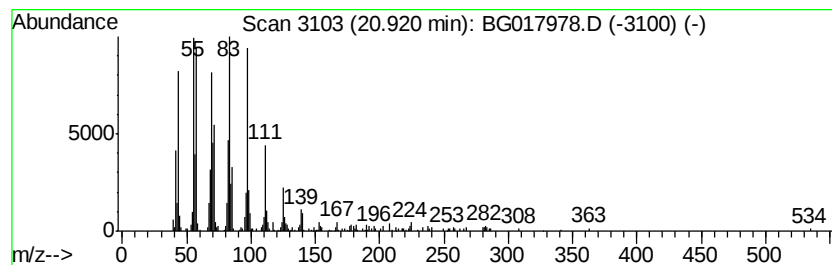
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG071715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 5-Eicosene, (E)- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.92	2.29 ng	203842	Chrysene-d12	21.20

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Eicosene, (E)-	280	C20H40	074685-30-6	97
2		1-Nonadecene	266	C19H38	018435-45-5	91
3		1-Docosene	308	C22H44	001599-67-3	86
4		Trichloroacetic acid, 4-hexadecy...	386	C18H33Cl3O2	1000282-92-4	83
5		1-Octadecanol	270	C18H38O	000112-92-5	81



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG071915\
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 Acq On : 20 Jul 2015 3:37
 Operator : TP/UM
 Sample : G3009-01
 Misc :
 ALS Vial : 54 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 MW-01-071515

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG071715.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...	2.86	58.5	ng	1593490	1	7.58	544285	20.0
2-Pentanone, 4-hy...	4.73	3.2	ng	88237	1	7.58	544285	20.0
unknown7.12	7.12	74.6	ng	2031060	1	7.58	544285	20.0
Benzene, 2-ethyl-...	8.68	3.7	ng	99639	1	7.58	544285	20.0
Benzene, 1,2,4,5-...	9.77	4.3	ng	196916	2	10.36	905512	20.0
Naphthalene, 1,2,...	10.83	2.4	ng	106172	2	10.36	905512	20.0
Naphthalene, 1,2,...	10.94	2.2	ng	101271	2	10.36	905512	20.0
Naphthalene, 1,2,...	11.92	3.0	ng	136823	2	10.36	905512	20.0
Dodecanoic acid	14.70	11.3	ng	739537	3	14.24	1312180	20.0
(Benzo(b)thien-6-...	14.85	2.7	ng	178719	3	14.24	1312180	20.0
Benzophenone	15.61	3.6	ng	238173	3	14.24	1312180	20.0
n-Hexadecanoic acid	17.91	3.3	ng	260160	4	17.00	1555760	20.0
5-Eicosene, (E)-	20.92	2.3	ng	203842	5	21.20	1776980	20.0