

Data Path : Z:\HPCHEM1\BNA G\DATA\BG050515\
 Data File : BG016880.D
 Acq On : 6 May 2015 1:29
 Operator : TP/IZ
 Sample : G2114-18
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 GRID-6(12-18)

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-BG050515.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	2.789	13	17	35	rBV	6155555	8328179	100.00%	19.337%
2	2.924	36	40	49	rVB2	171788	314507	3.78%	0.730%
3	3.007	49	54	59	rBV	218036	281474	3.38%	0.654%
4	4.922	374	380	387	rBV	182353	265665	3.19%	0.617%
5	5.386	452	459	470	rBV	1635098	2385402	28.64%	5.539%
6	6.972	720	729	740	rVB	1739115	2863724	34.39%	6.649%
7	7.337	784	791	803	rBV	1681424	2606467	31.30%	6.052%
8	7.801	863	870	877	rVB	312707	508765	6.11%	1.181%
9	8.124	918	925	934	rVB	1073888	1749262	21.00%	4.062%
10	8.958	1059	1067	1076	rBV	1019514	1670235	20.06%	3.878%
11	10.598	1338	1346	1354	rBV	436974	752254	9.03%	1.747%
12	12.801	1715	1721	1727	rBV	91867	136318	1.64%	0.317%
13	13.065	1758	1766	1772	rBV	2511320	3748808	45.01%	8.704%
14	13.894	1902	1907	1918	rBV	102019	156166	1.88%	0.363%
15	14.440	1994	2000	2010	rBV2	681944	1039079	12.48%	2.413%
16	15.803	2225	2232	2237	rVB	182049	245272	2.95%	0.569%
17	15.927	2245	2253	2269	rBV	2154795	3028318	36.36%	7.031%
18	17.184	2461	2467	2473	rVB2	913172	1275189	15.31%	2.961%
19	18.089	2613	2621	2630	rBV2	397990	607212	7.29%	1.410%
20	19.370	2834	2839	2846	rBV	194563	295957	3.55%	0.687%
21	19.816	2909	2915	2920	rBV	4427306	5482332	65.83%	12.729%
22	20.504	3027	3032	3038	rBV2	94528	181000	2.17%	0.420%
23	21.079	3126	3130	3140	rBV	1205847	1397140	16.78%	3.244%
24	21.361	3172	3178	3189	rBV	1151505	1441308	17.31%	3.347%
25	21.984	3279	3284	3291	rBV2	73694	117556	1.41%	0.273%
26	22.184	3314	3318	3324	rVB	71502	98672	1.18%	0.229%
27	22.584	3382	3386	3393	rVB	176203	251542	3.02%	0.584%
28	22.971	3447	3452	3456	rBV4	43241	92132	1.11%	0.214%
29	23.671	3563	3571	3580	rVB2	682391	1348970	16.20%	3.132%
30	24.346	3681	3686	3699	rVB4	135293	315664	3.79%	0.733%
31	26.144	3988	3992	4002	rVB4	35633	83453	1.00%	0.194%

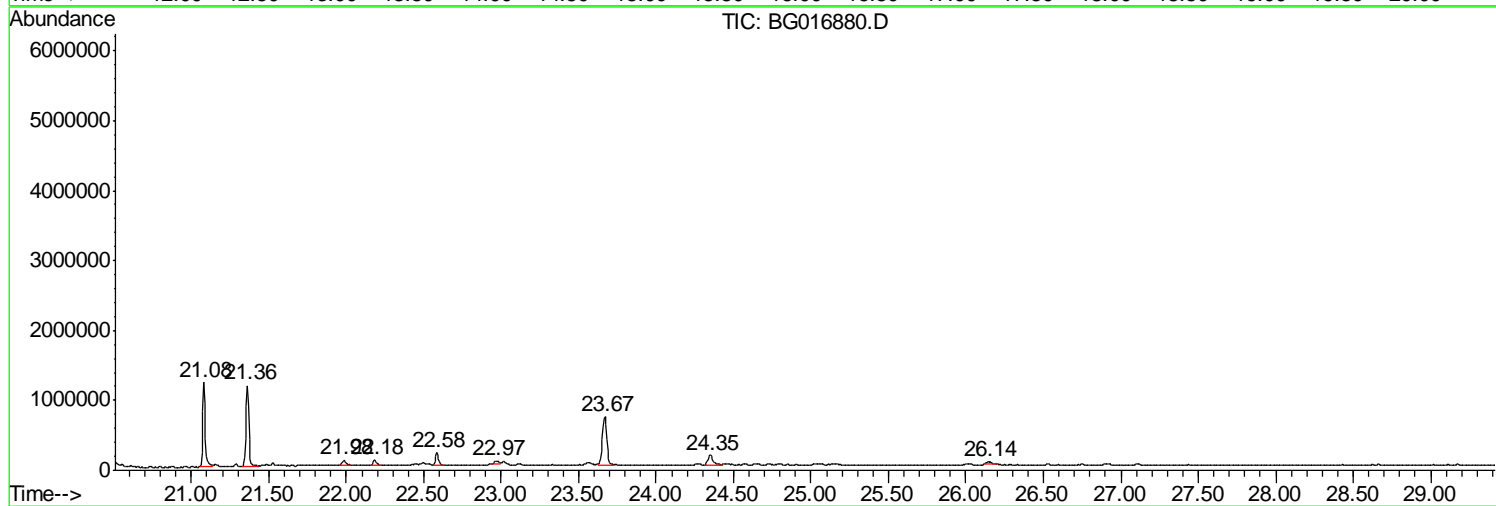
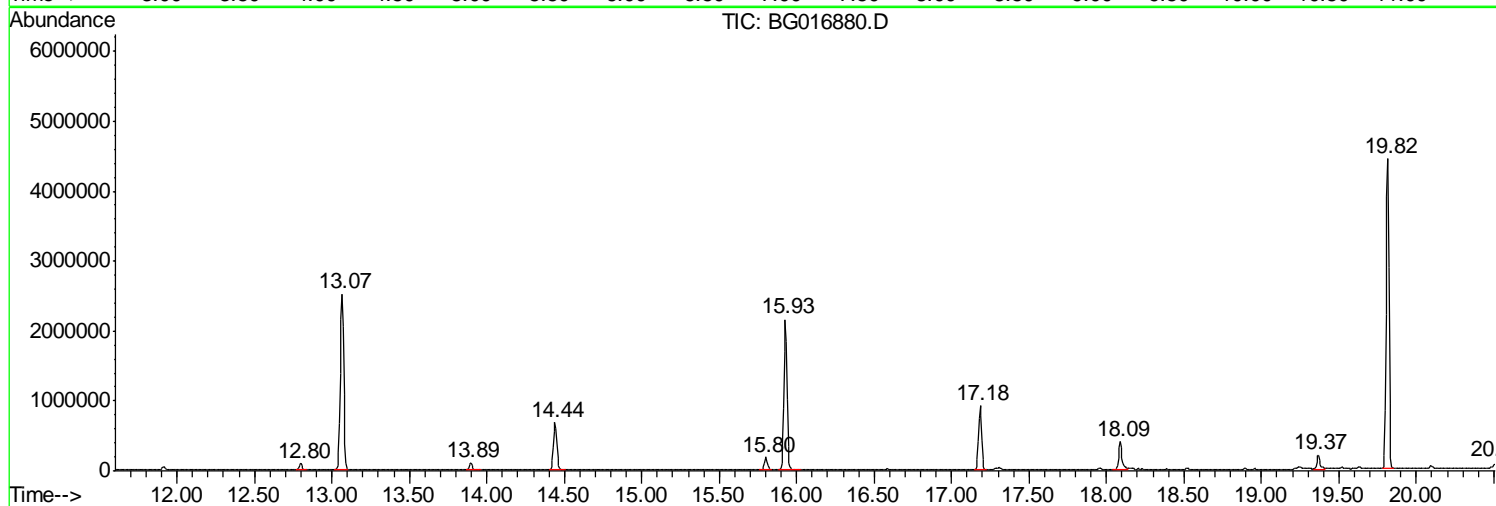
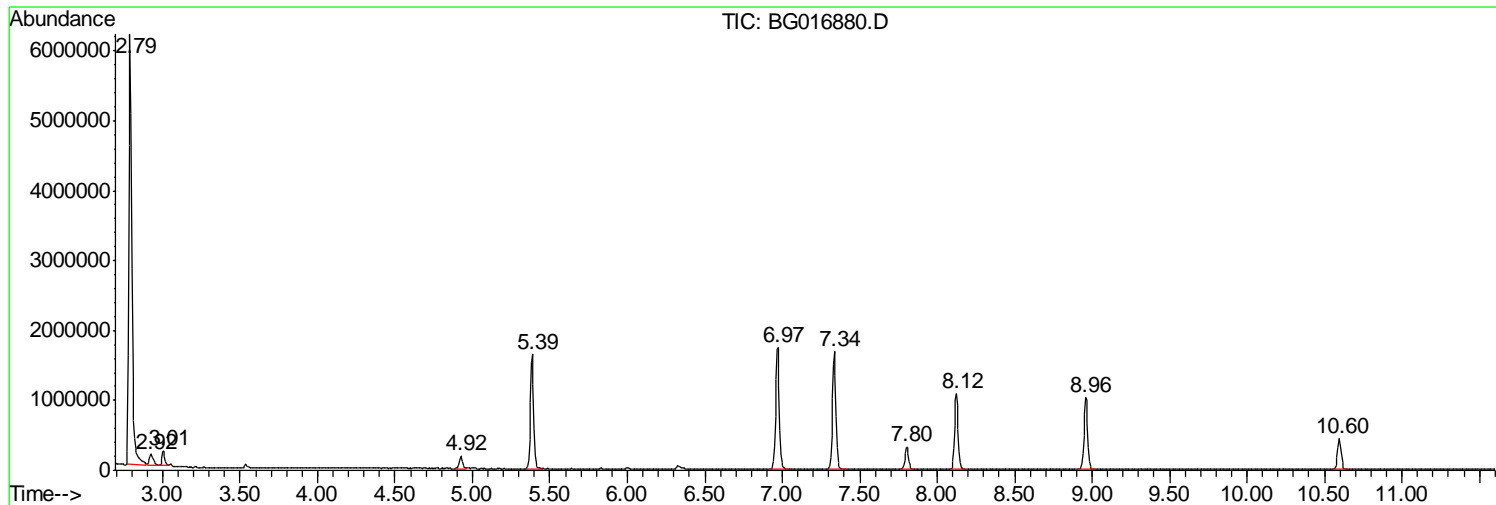
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Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG050515.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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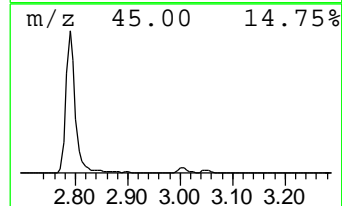
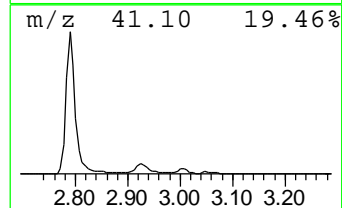
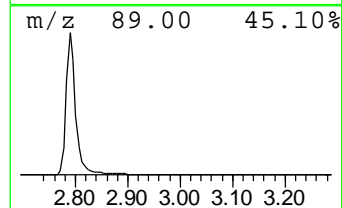
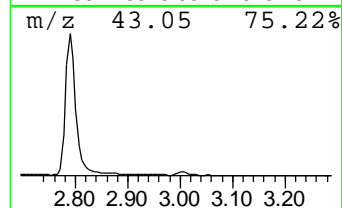
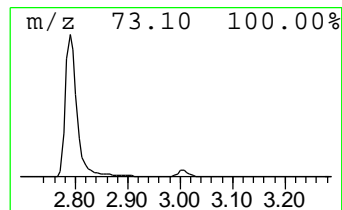
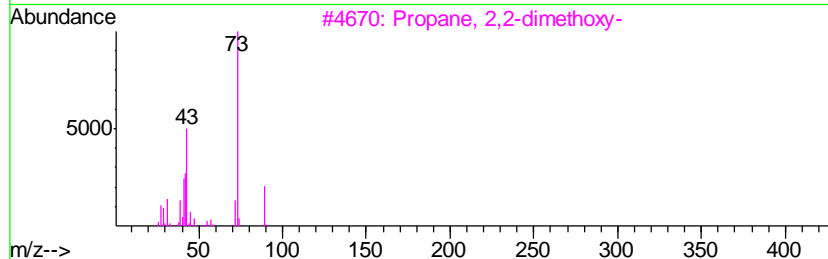
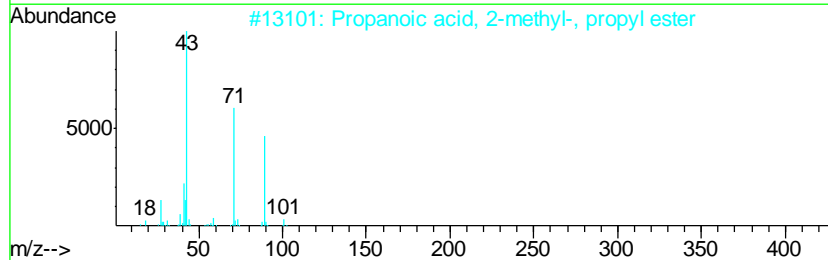
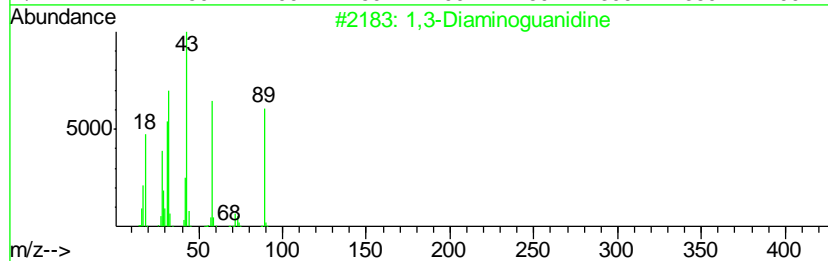
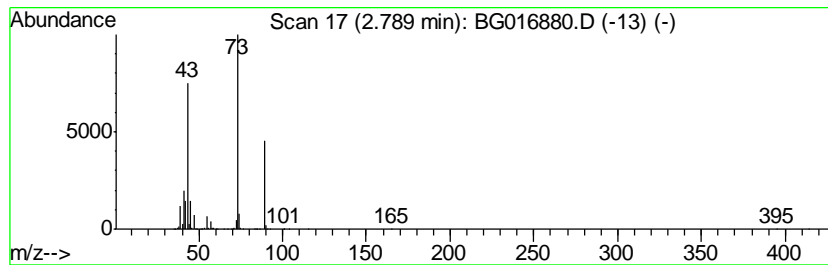
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TIC Library : C:\DATABASE\NIST02.L
 TIC Integration Parameters: LSCINT.P

 Peak Number 1 1,3-Diaminoguanidine Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.79	327.39 ng	8328180	1,4-Dichlorobenzene-d4	7.80

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,3-Diaminoguanidine	89	CH7N5	004364-78-7	53
2		Propanoic acid, 2-methyl-, propy...	130	C7H14O2	000644-49-5	47
3		Propane, 2,2-dimethoxy-	104	C5H12O2	000077-76-9	42
4		Acetamide, N-methyl-	73	C3H7NO	000079-16-3	25
5		2-Butanone, 3-methoxy-3-methyl-	116	C6H12O2	036687-98-6	25



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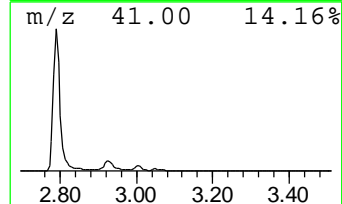
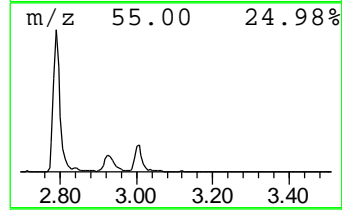
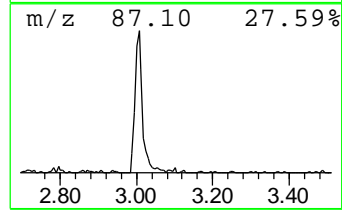
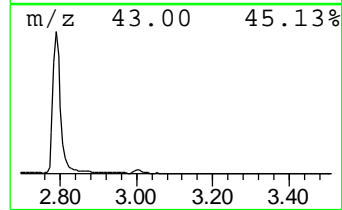
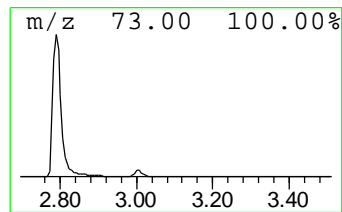
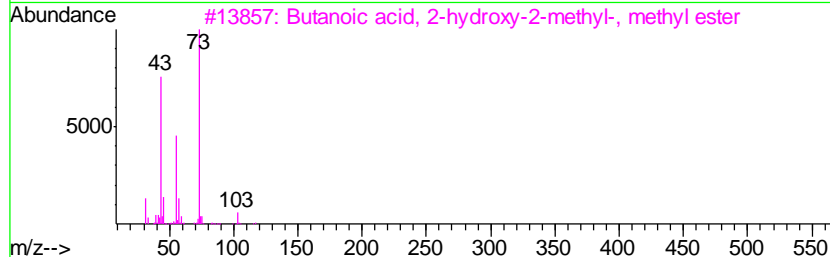
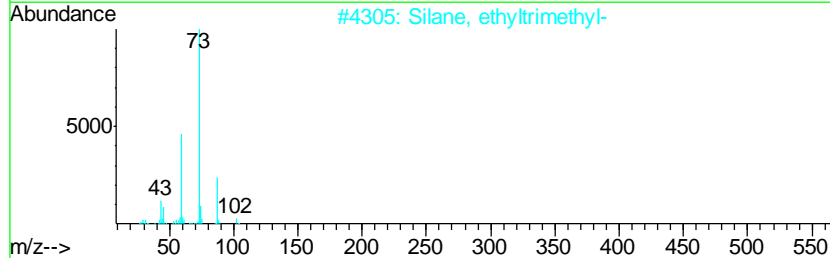
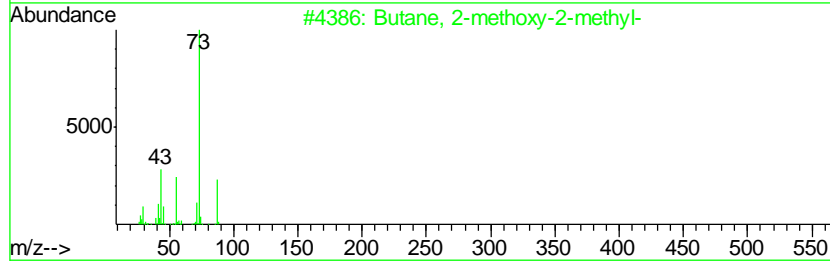
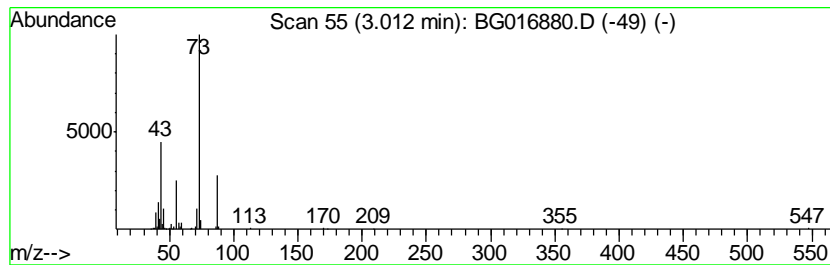
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 Peak Number 3 Butane, 2-methoxy-2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.01	11.07 ng	281474	1,4-Dichlorobenzene-d4	7.80

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2		Silane, ethyltrimethyl-	102	C5H14Si	003439-38-1	36
3		Butanoic acid, 2-hydroxy-2-methyl...	132	C6H12O3	032793-34-3	10
4		Methane, isothiocyanato-	73	C2H3NS	000556-61-6	9
5		Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	9



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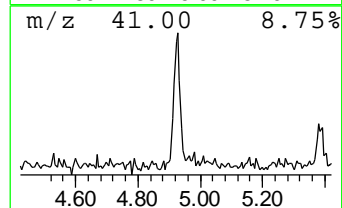
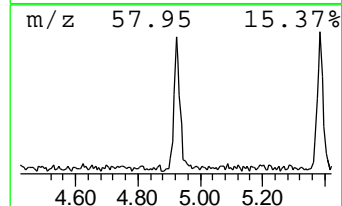
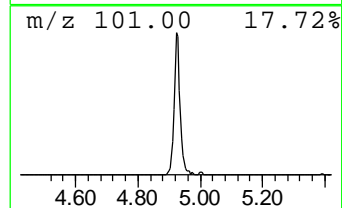
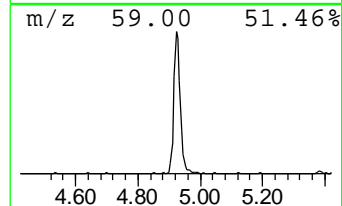
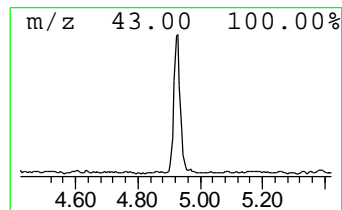
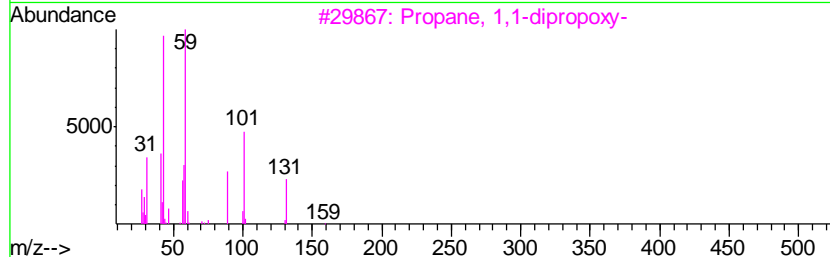
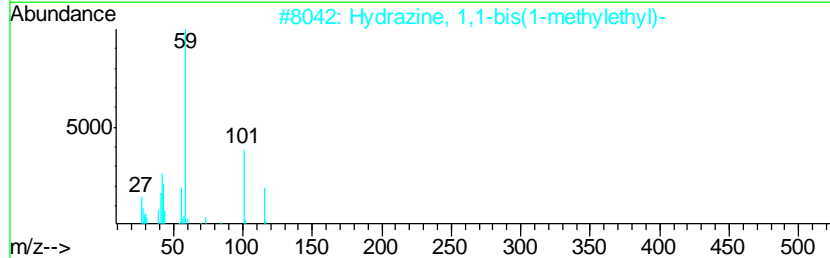
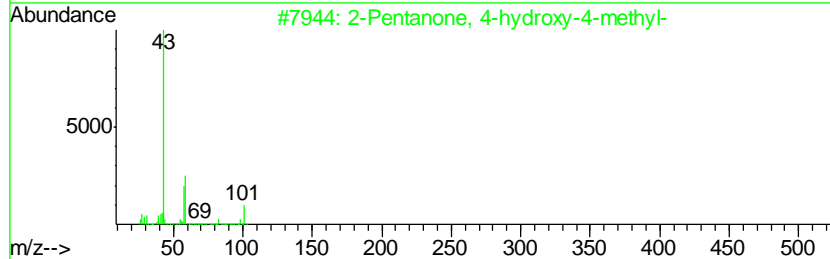
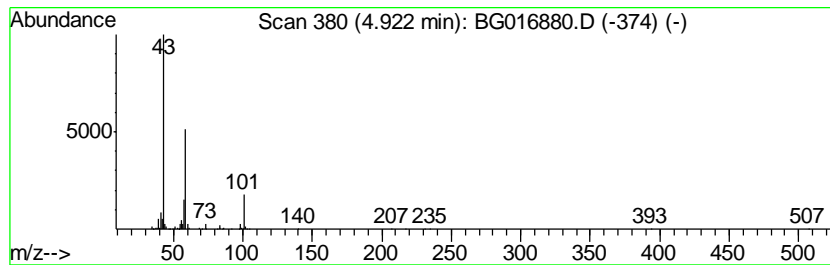
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TIC Library : C:\DATABASE\NIST02.L
 TIC Integration Parameters: LSCINT.P

 Peak Number 4 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.92	10.44 ng	265665	1,4-Dichlorobenzene-d4	7.80

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	53
2		Hydrazine, 1,1-bis(1-methylethyl)-	116	C6H16N2	000921-14-2	37
3		Propane, 1,1-dipropoxy-	160	C9H20O2	004744-11-0	36
4		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	35
5		Propane, 1-ethoxy-2-methyl-	102	C6H14O	000627-02-1	25



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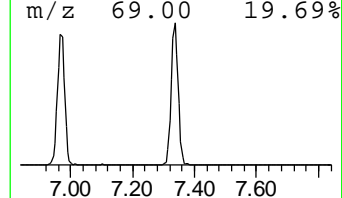
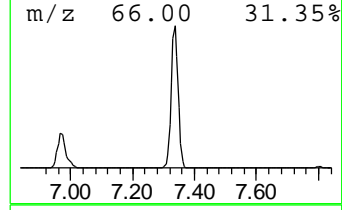
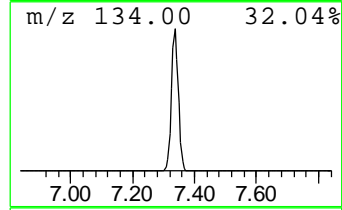
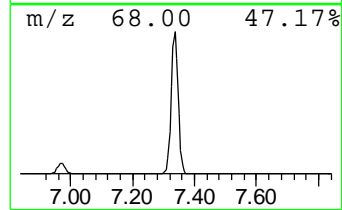
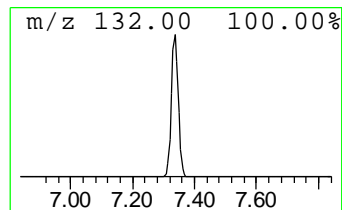
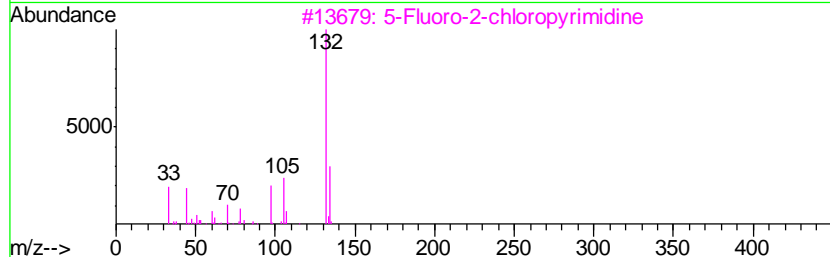
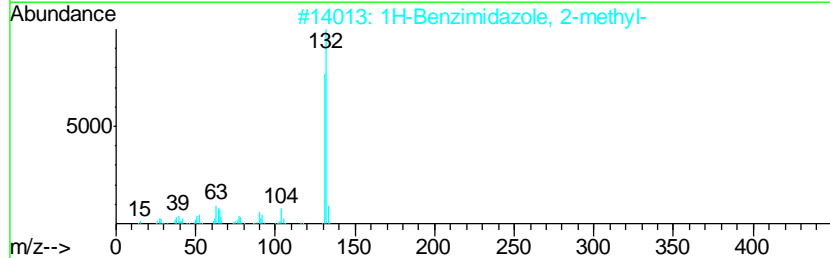
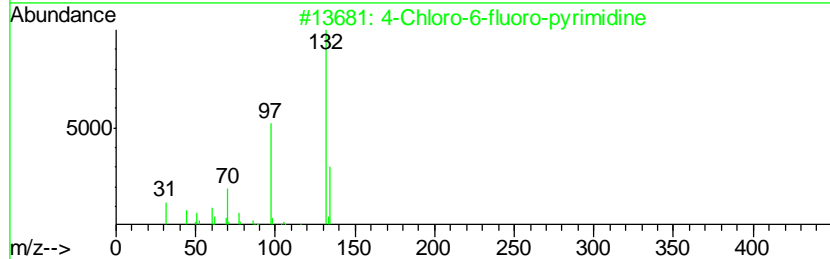
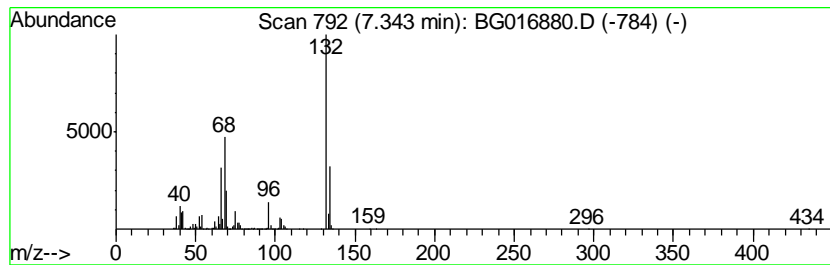
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TIC Library : C:\DATABASE\NIST02.L
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 Peak Number 5 unknown7.34 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.34	102.46 ng	2606470	1,4-Dichlorobenzene-d4	7.80

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	27
2		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
3		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
4		Pyrazolo(2,3-a)pyridine, 7-methyl-	132	C8H8N2	034760-58-2	12
5		Benzene, 1-methyl-4-(1-methyleth...	132	C10H12	001195-32-0	10



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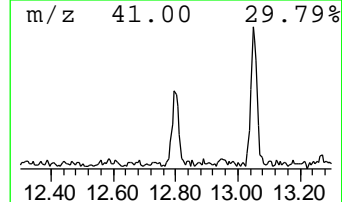
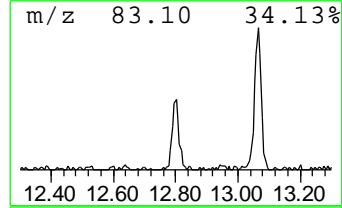
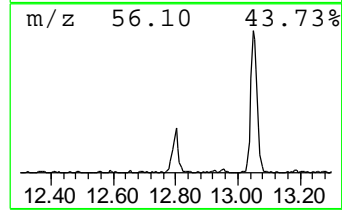
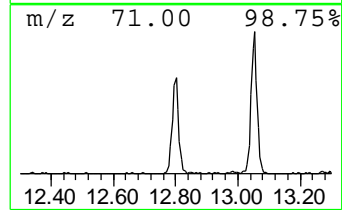
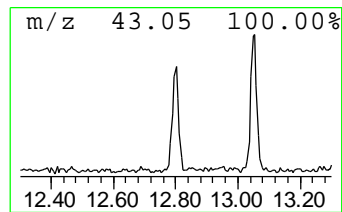
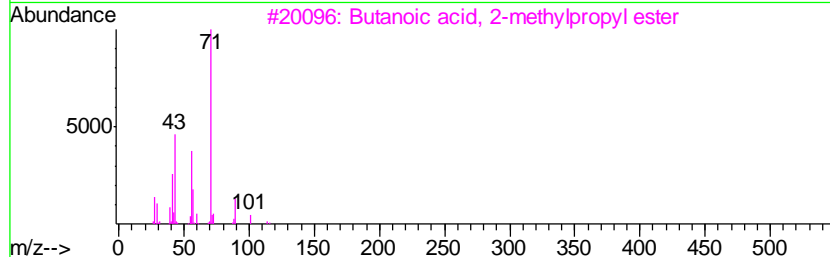
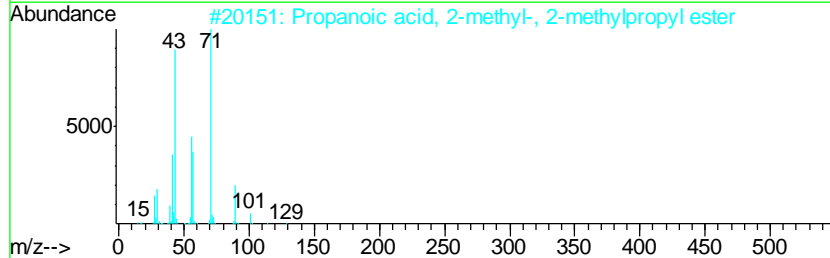
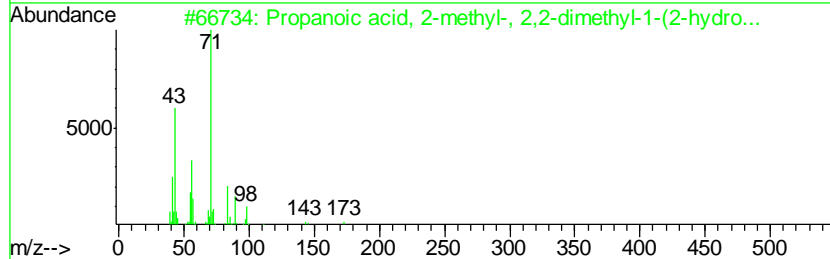
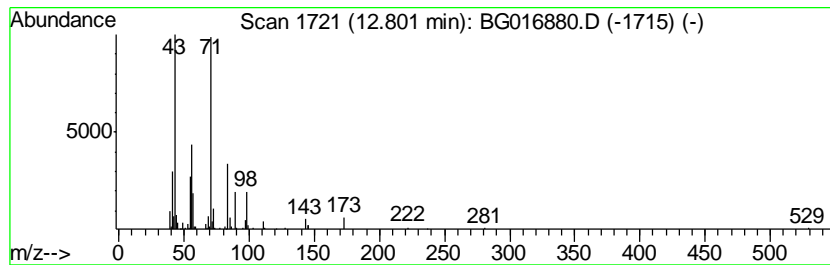
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 Peak Number 7 Propanoic acid, 2-methyl-, ... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.80	2.62 ng	136318	Acenaphthene-d10	14.44

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Propanoic acid, 2-methyl-, 2,2-d...	216	C12H24O3	074367-33-2	56
2		Propanoic acid, 2-methyl-, 2-met...	144	C8H16O2	000097-85-8	53
3		Butanoic acid, 2-methylpropyl ester	144	C8H16O2	000539-90-2	53
4		2,2,4-Trimethyl-1,3-pentanediol ...	286	C16H30O4	006846-50-0	50
5		Butanoic acid, 1-methylpropyl ester	144	C8H16O2	000819-97-6	47



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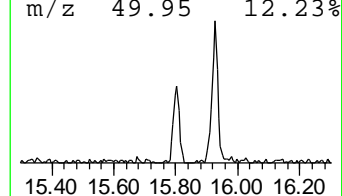
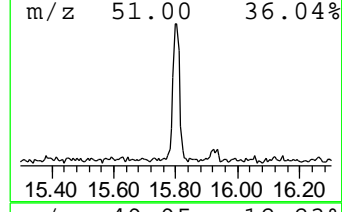
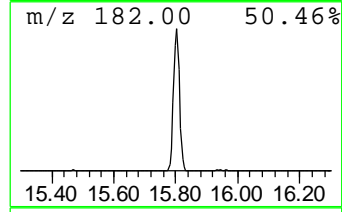
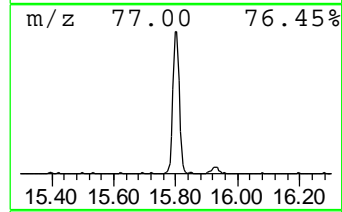
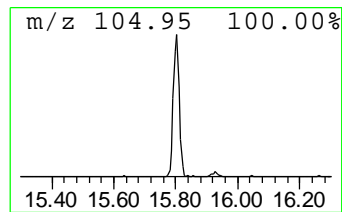
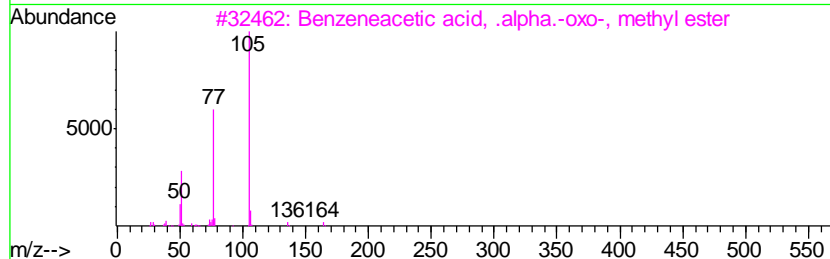
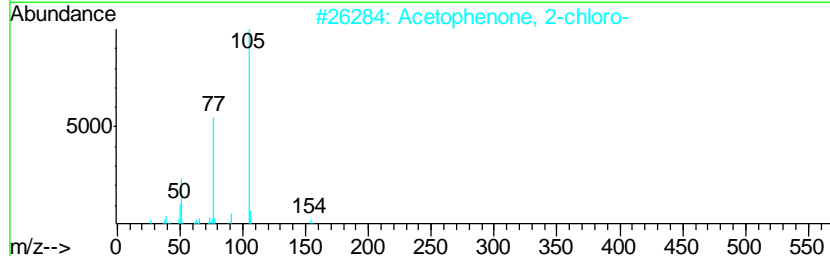
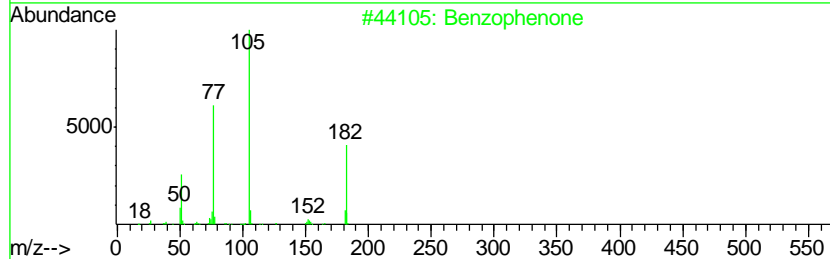
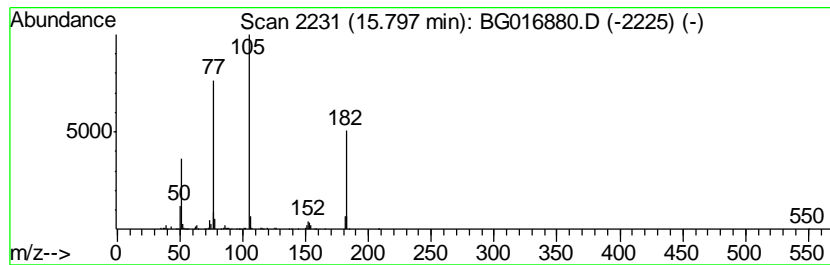
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 8 Benzophenone Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.80	4.72 ng	245272	Acenaphthene-d10	14.44

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzophenone	182	C13H10O	000119-61-9	97
2		Acetophenone, 2-chloro-	154	C8H7ClO	000532-27-4	49
3		Benzeneacetic acid, .alpha.-oxo-...	164	C9H8O3	015206-55-0	49
4		Benzoylformic acid	150	C8H6O3	000611-73-4	49
5		Glyoxylohydroximoyl chloride, 2-...	183	C8H6ClNO2	004937-87-5	47



Data Path : Z:\HPCHEM1\BNA G\DATA\BG050515\
 Data File : BG016880.D
 Acq On : 6 May 2015 1:29
 Operator : TP/IZ
 Sample : G2114-18
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampled :
 GRID-6(12-18)

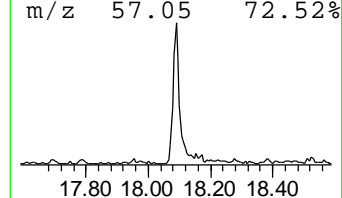
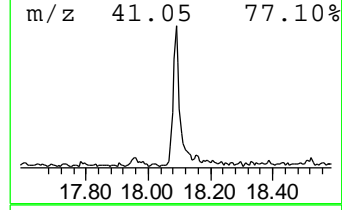
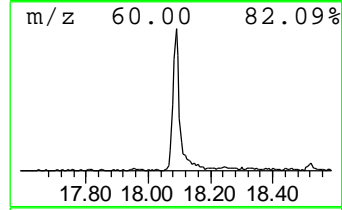
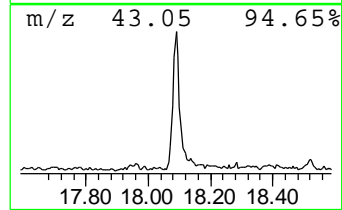
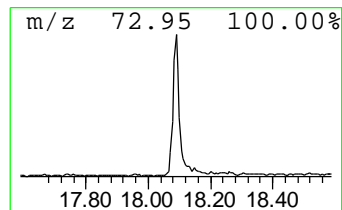
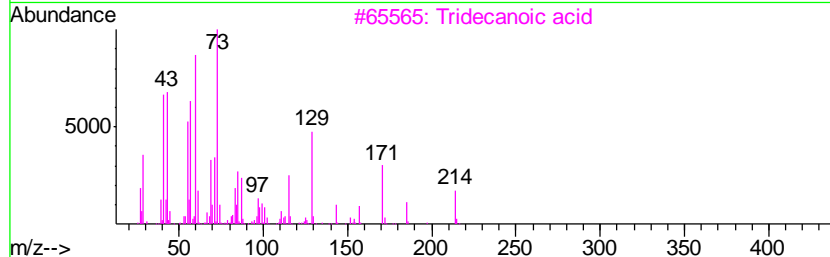
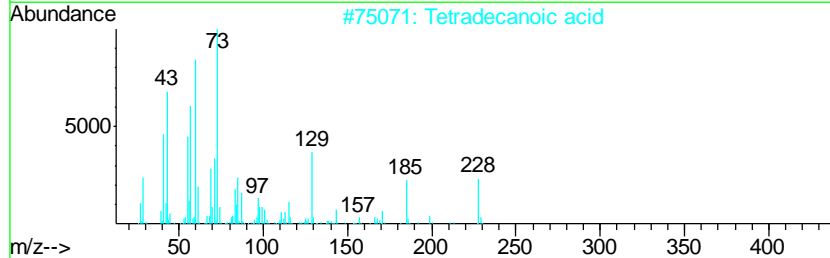
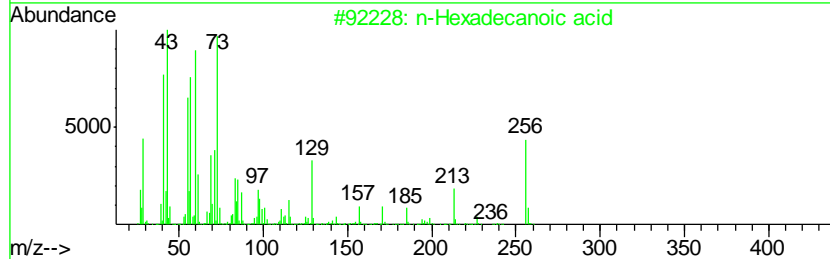
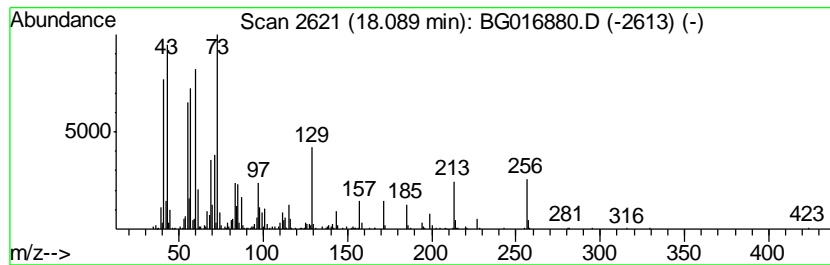
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 9 n-Hexadecanoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.09	9.52 ng	607212	Phenanthrene-d10	17.18

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2		Tetradecanoic acid	228	C14H28O2	000544-63-8	86
3		Tridecanoic acid	214	C13H26O2	000638-53-9	64
4		n-Decanoic acid	172	C10H20O2	000334-48-5	60
5		Pentadecanoic acid	242	C15H30O2	001002-84-2	53



Data Path : Z:\HPCHEM1\BNA G\DATA\BG050515\
 Data File : BG016880.D
 Acq On : 6 May 2015 1:29
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 Misc :
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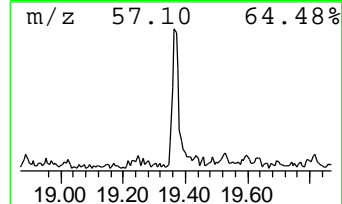
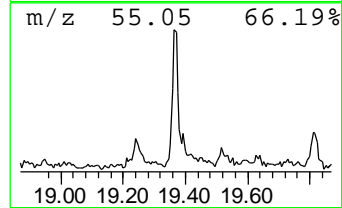
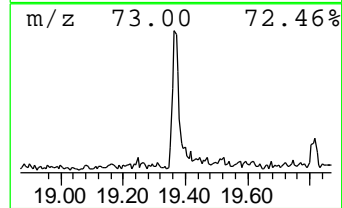
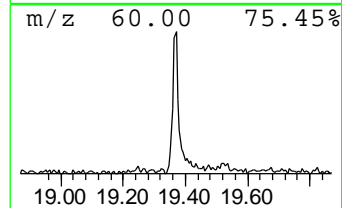
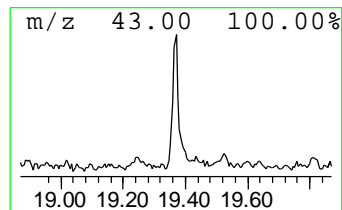
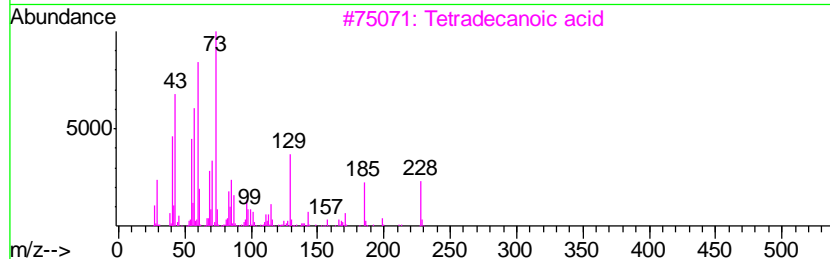
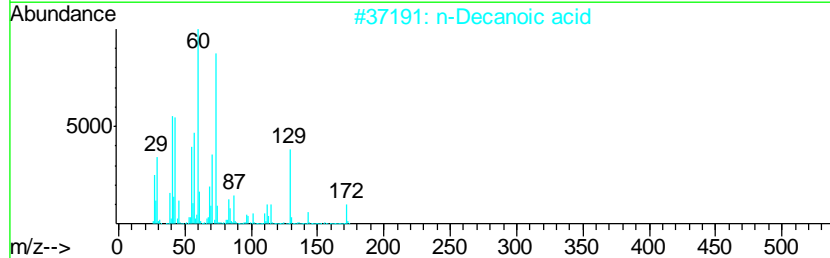
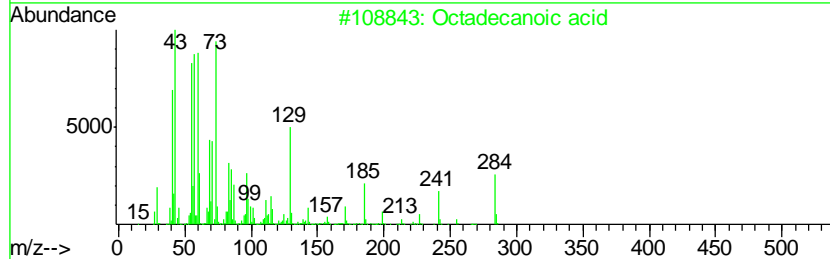
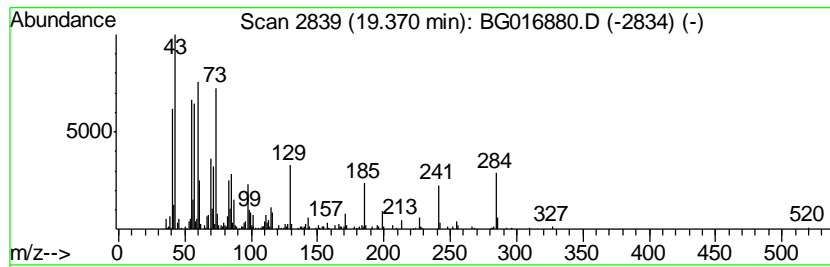
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 10 Octadecanoic acid Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.37	4.11 ng	295957	Chrysene-d12	21.36

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecanoic acid	284	C18H36O2	000057-11-4	97
2		n-Decanoic acid	172	C10H20O2	000334-48-5	50
3		Tetradecanoic acid	228	C14H28O2	000544-63-8	49
4		Undecanoic acid	186	C11H22O2	000112-37-8	45
5		Dodecanoic acid	200	C12H24O2	000143-07-7	43



Data Path : Z:\HPCHEM1\BNA G\DATA\BG050515\
 Data File : BG016880.D
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 Misc :
 ALS Vial : 19 Sample Multiplier: 1

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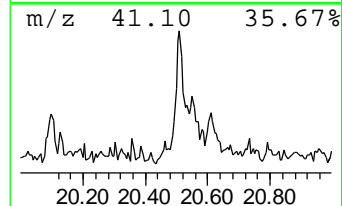
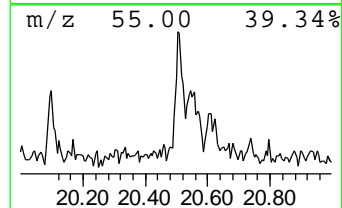
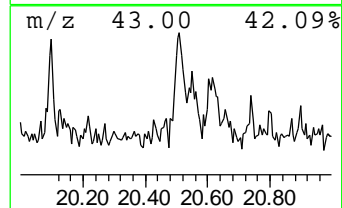
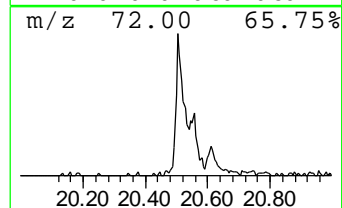
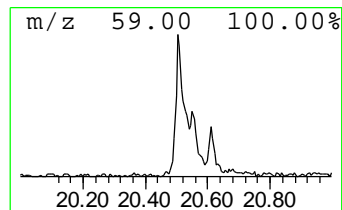
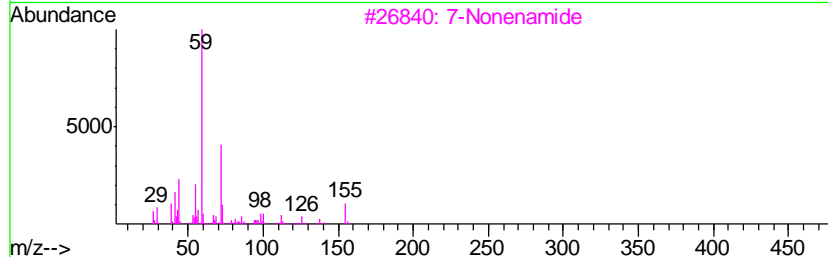
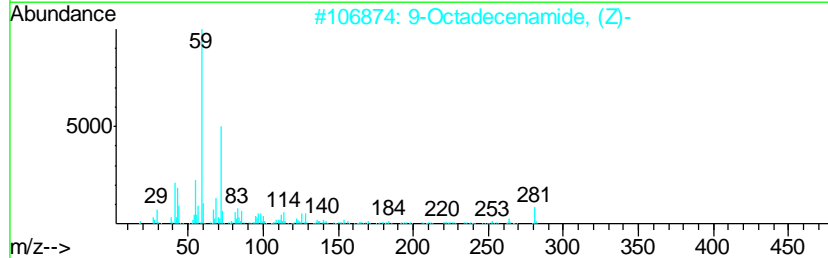
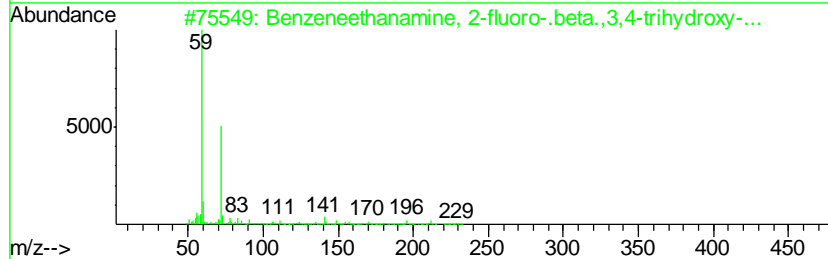
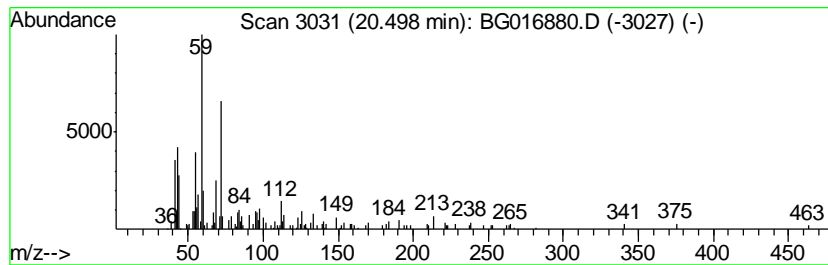
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 11 Benzeneethanamine, 2-fluoro... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.50	2.51 ng	181000	Chrysene-d12	21.36

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzeneethanamine, 2-fluoro-.bet...	229	C11H16FNO3	061338-98-5	72
2		9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	68
3		7-Nonenamide	155	C9H17NO	090949-53-4	53
4		Hexadecanamide	255	C16H33NO	000629-54-9	50
5		2,2-Dimethyl-4-hydroxymethyl-5-(...	198	C11H18O3	1000284-41-5	38



Data Path : Z:\HPCHEM1\BNA G\DATA\BG050515\
 Data File : BG016880.D
 Acq On : 6 May 2015 1:29
 Operator : TP/IZ
 Sample : G2114-18
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

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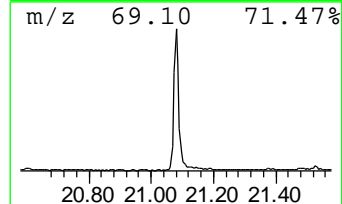
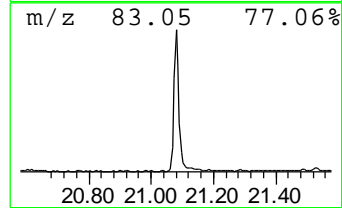
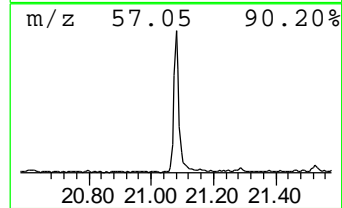
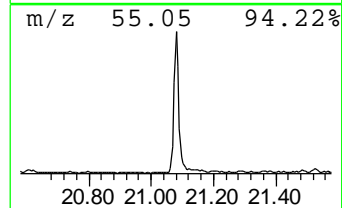
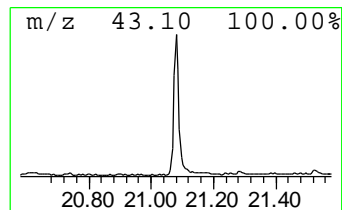
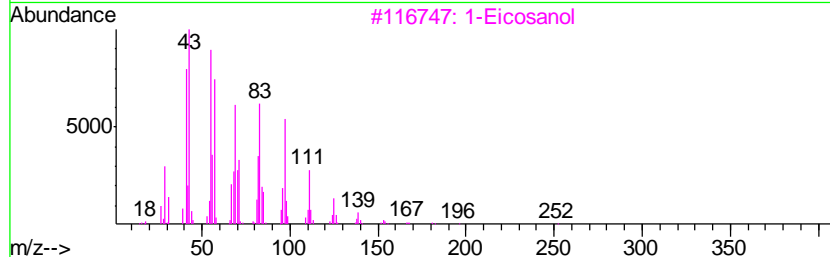
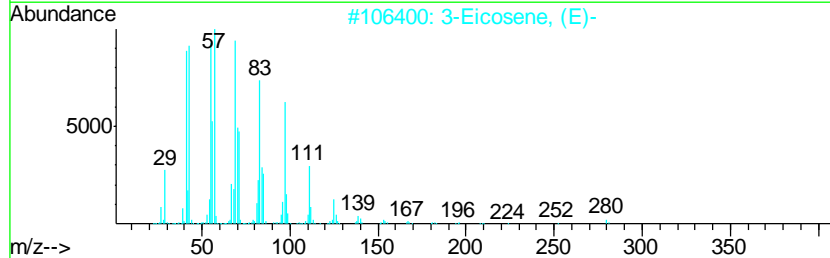
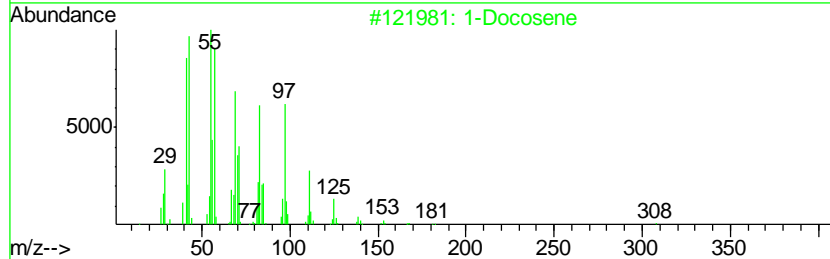
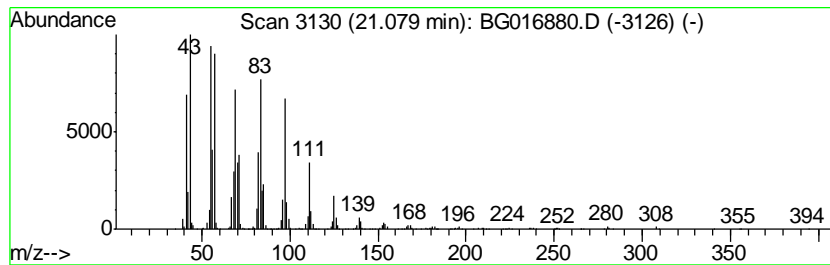
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 12 1-Docosene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.08	19.39 ng	1397140	Chrysene-d12	21.36

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Docosene	308	C22H44	001599-67-3	99
2		3-Eicosene, (E)-	280	C20H40	074685-33-9	96
3		1-Eicosanol	298	C20H42O	000629-96-9	94
4		1-Octadecanol	270	C18H38O	000112-92-5	91
5		1-Docosanol	326	C22H46O	000661-19-8	91



Data Path : Z:\HPCHEM1\BNA G\DATA\BG050515\
 Data File : BG016880.D
 Acq On : 6 May 2015 1:29
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 Misc :
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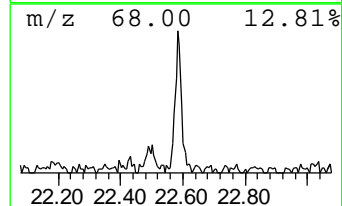
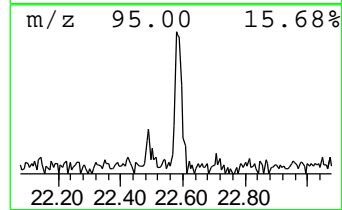
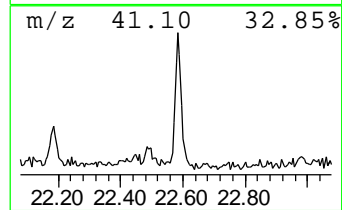
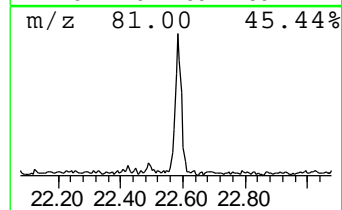
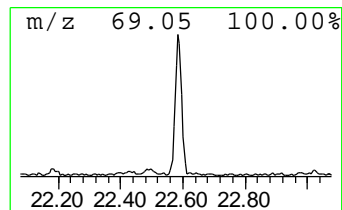
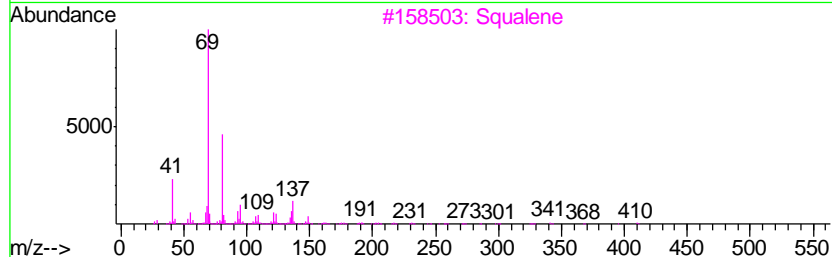
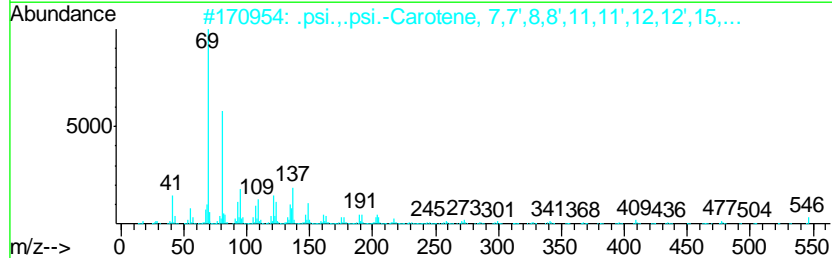
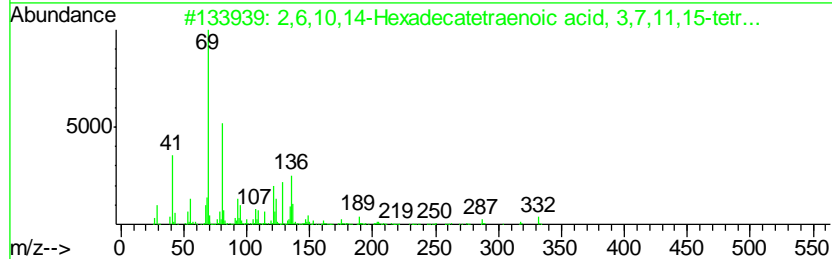
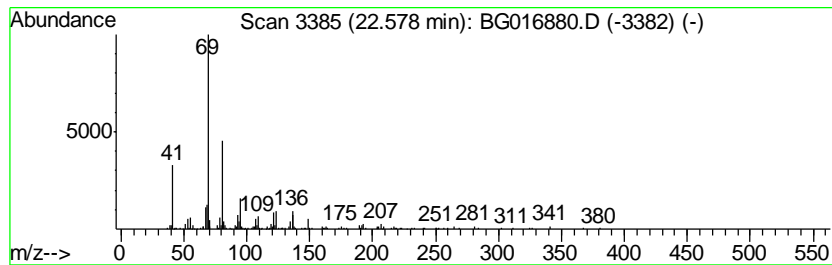
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 13 2,6,10,14-Hexadecatetraenoic... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.58	3.73 ng	251542	Perylene-d12	23.66

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2,6,10,14-Hexadecatetraenoic aci...	332	C22H36O2	024035-35-6	72
2		.psi.,.psi.-Carotene, 7,7',8,8',...	547	C40H66	000502-62-5	72
3		Squalene	410	C30H50	007683-64-9	72
4		7,11-Dimethyldodeca-2,6,10-trien...	208	C14H24O	125529-10-4	64
5		2,6,10-Dodecatrien-1-ol, 3,7,11-...	222	C15H26O	004602-84-0	59



Data Path : Z:\HPCHEM1\BNA G\DATA\BG050515\
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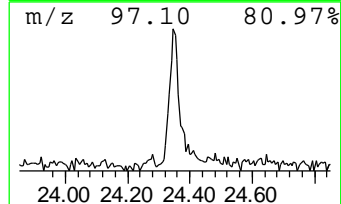
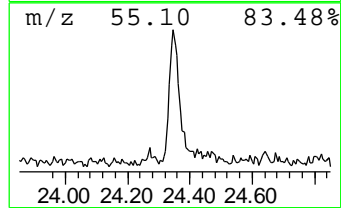
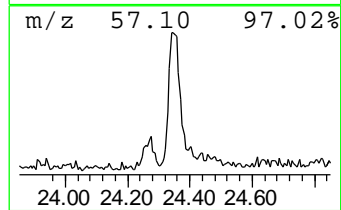
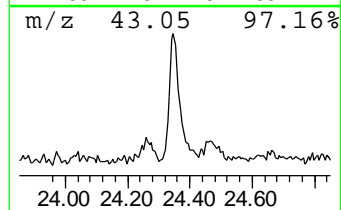
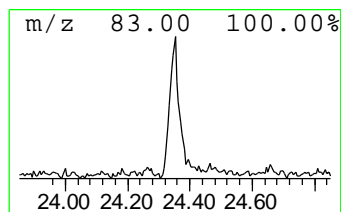
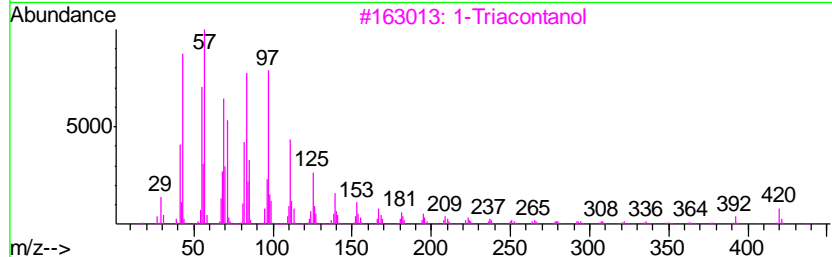
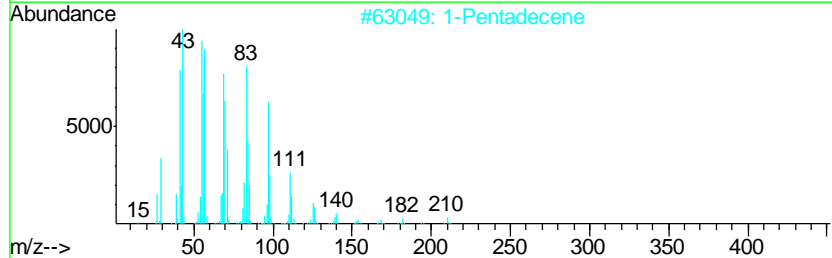
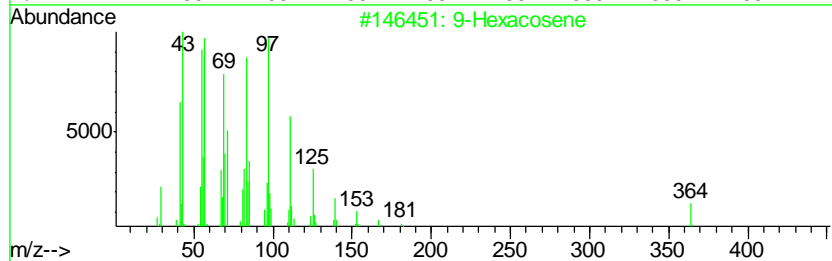
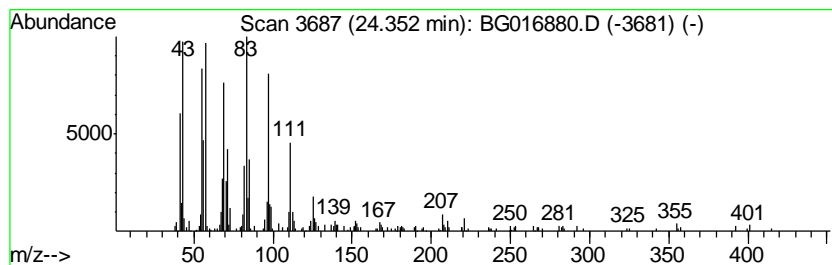
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 14 9-Hexacosene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.35	4.68 ng	315664	Perylene-d12	23.66

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	9-Hexacosene	364	C26H52	071502-22-2	89
2		1-Pentadecene	210	C15H30	013360-61-7	81
3		1-Triacontanol	438	C30H62O	000593-50-0	64
4		1-Tetracosanol	354	C24H50O	000506-51-4	64
5		1-Docosene	308	C22H44	001599-67-3	64



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG050515\
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 Acq On : 6 May 2015 1:29
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 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
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Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG050515.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
1,3-Diaminoguanidine	2.79	327.4	ng	8328180	1	7.80	508765	20.0
Butane, 2-methoxy...	3.01	11.1	ng	281474	1	7.80	508765	20.0
2-Pentanone, 4-hy...	4.92	10.4	ng	265665	1	7.80	508765	20.0
unknown7.34	7.34	102.5	ng	2606470	1	7.80	508765	20.0
Propanoic acid, 2...	12.80	2.6	ng	136318	3	14.44	1039080	20.0
Benzophenone	15.80	4.7	ng	245272	3	14.44	1039080	20.0
n-Hexadecanoic acid	18.09	9.5	ng	607212	4	17.18	1275190	20.0
Octadecanoic acid	19.37	4.1	ng	295957	5	21.36	1441310	20.0
Benzeneethanamine...	20.50	2.5	ng	181000	5	21.36	1441310	20.0
1-Docosene	21.08	19.4	ng	1397140	5	21.36	1441310	20.0
2,6,10,14-Hexadec...	22.58	3.7	ng	251542	6	23.66	1348970	20.0
9-Hexacosene	24.35	4.7	ng	315664	6	23.66	1348970	20.0