

Data Path : Z:\HPCHEM1\BNA G\DATA\BG042715\
 Data File : BG016763.D
 Acq On : 28 Apr 2015 10:56
 Operator : TP/IZ
 Sample : G1991-11
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SB-74D

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-BG042315.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	2.723	3	6	11	rVB	75381	94670	1.14%	0.338%
2	4.680	331	339	349	rVB	137062	192484	2.32%	0.687%
3	5.168	416	422	437	rBV	766978	1089050	13.13%	3.889%
4	6.777	689	696	709	rBV	741251	1190885	14.36%	4.252%
5	7.112	746	753	763	rVB	841249	1259744	15.19%	4.498%
6	7.571	824	831	838	rVB	196871	299681	3.61%	1.070%
7	7.888	876	885	896	rBV	633973	979848	11.81%	3.499%
8	8.734	1020	1029	1044	rBV	436281	720804	8.69%	2.574%
9	10.356	1298	1305	1313	rBV	279733	455296	5.49%	1.626%
10	12.841	1721	1728	1738	rBV	1210543	1694439	20.42%	6.050%
11	14.228	1957	1964	1971	rBV2	415378	625770	7.54%	2.234%
12	15.591	2191	2196	2203	rBV	85631	114447	1.38%	0.409%
13	15.720	2212	2218	2226	rBV	790748	1089297	13.13%	3.890%
14	16.971	2422	2431	2436	rBV	480584	688638	8.30%	2.459%
15	17.876	2580	2585	2597	rBV5	64778	140908	1.70%	0.503%
16	18.076	2615	2619	2624	rVB	110670	151178	1.82%	0.540%
17	18.252	2644	2649	2655	rBV	397274	540649	6.52%	1.930%
18	18.440	2676	2681	2691	rVB	705320	939710	11.33%	3.355%
19	18.593	2701	2707	2712	rVB	543241	711762	8.58%	2.541%
20	19.098	2786	2793	2798	rBV	6299235	8295916	100.00%	29.622%
21	19.604	2874	2879	2885	rBV	1517307	1945663	23.45%	6.947%
22	19.750	2900	2904	2908	rBV	109532	137076	1.65%	0.489%
23	19.815	2908	2915	2921	rVB	1587587	2001389	24.12%	7.146%
24	20.027	2947	2951	2954	rBV	144096	171741	2.07%	0.613%
25	20.873	3091	3095	3104	rVB	115585	176189	2.12%	0.629%
26	21.161	3138	3144	3151	rBV	540118	739447	8.91%	2.640%
27	21.484	3195	3199	3204	rBV	128233	145121	1.75%	0.518%
28	21.748	3240	3244	3254	rBV	392823	589364	7.10%	2.104%
29	22.224	3321	3325	3330	rBV2	102736	139932	1.69%	0.500%
30	23.352	3511	3517	3530	rBV	338532	684584	8.25%	2.444%

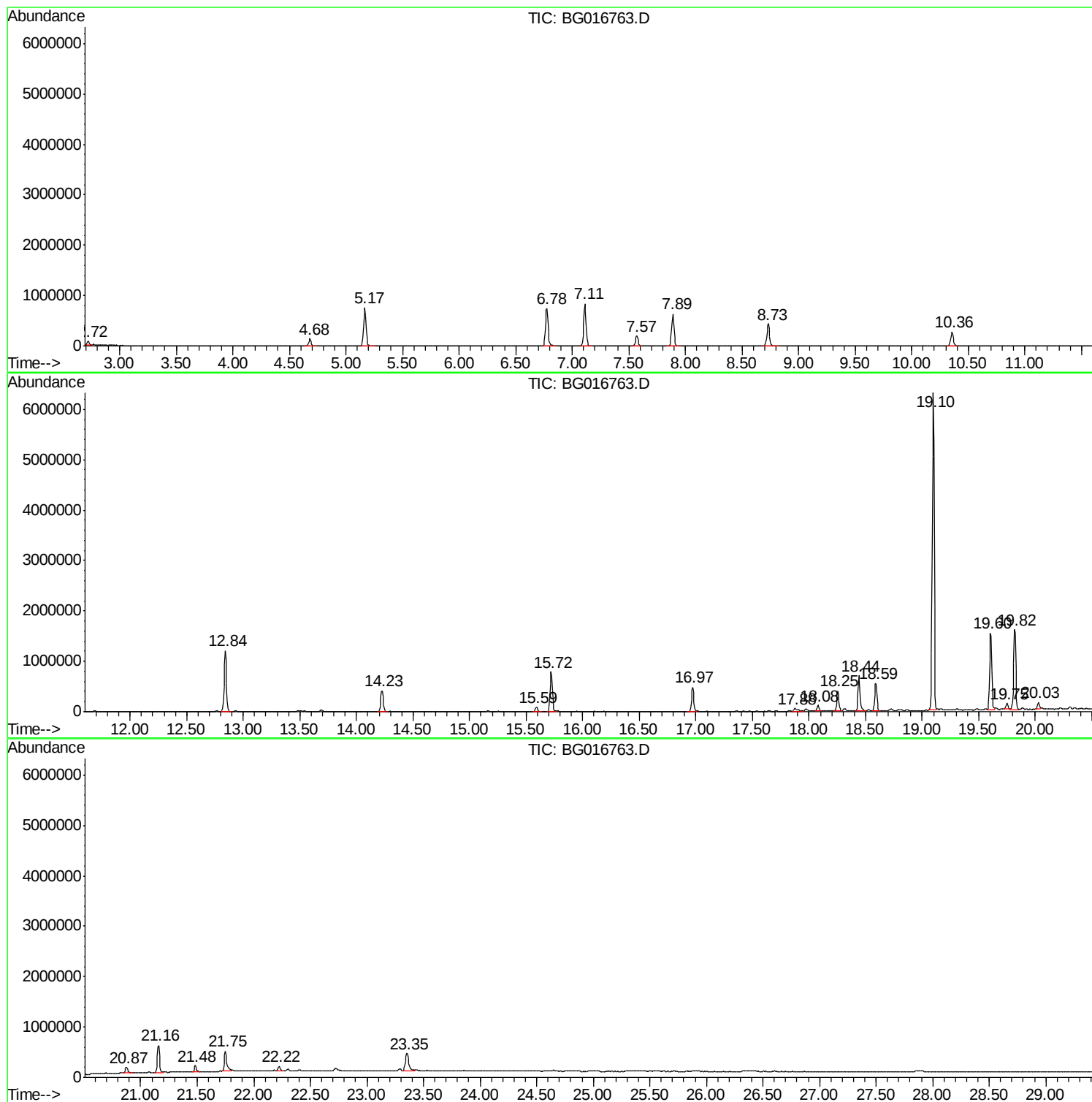
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Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG042315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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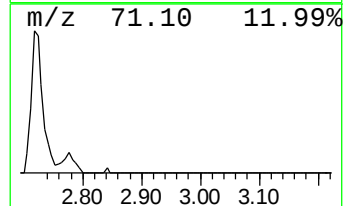
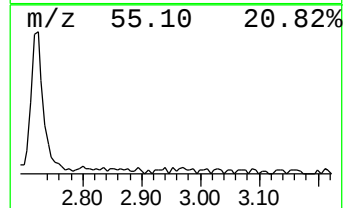
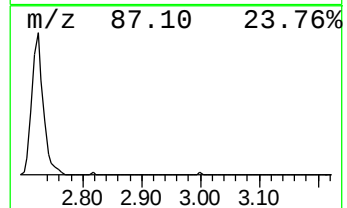
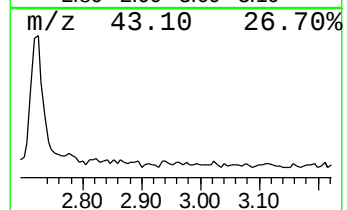
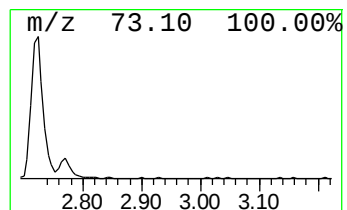
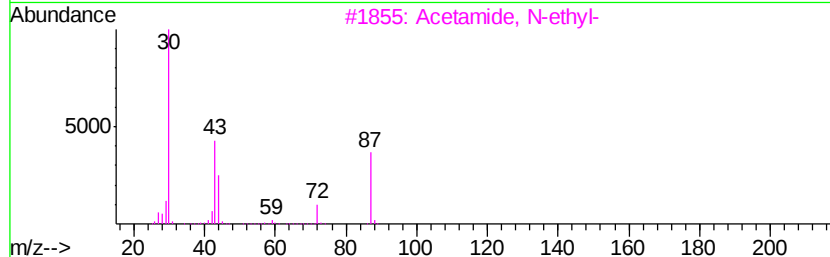
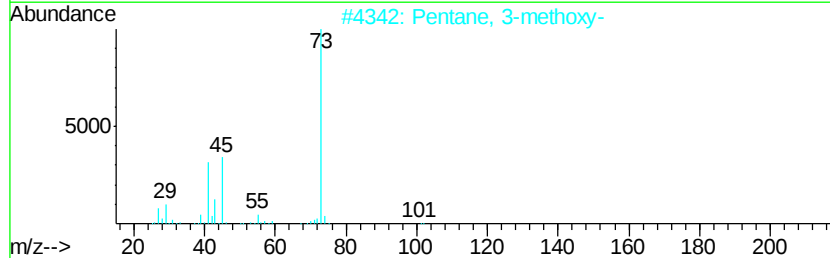
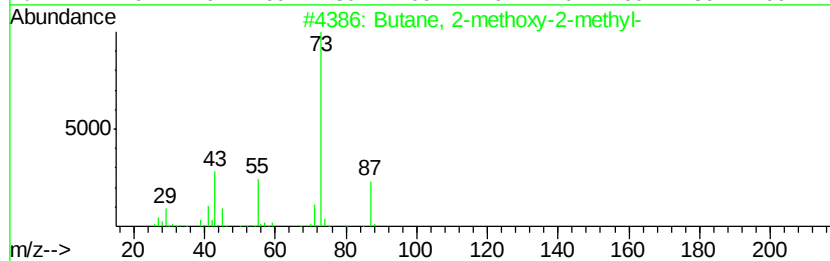
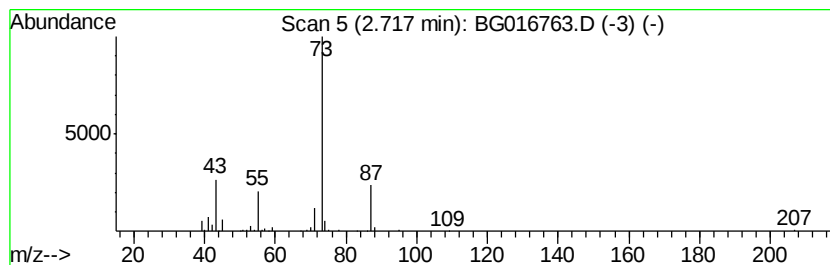
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.72	6.32 ng	94670	1,4-Dichlorobenzene-d4	7.57

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	72
2			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	17
3			Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	9
4			Silane, tetramethyl-	88	C4H12Si	000075-76-3	9
5			1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	9



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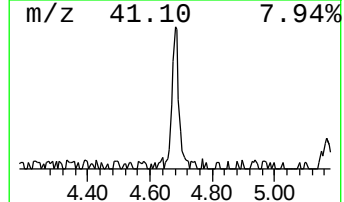
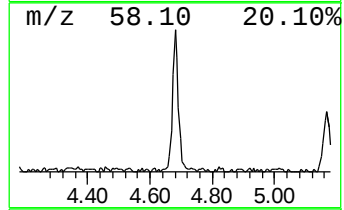
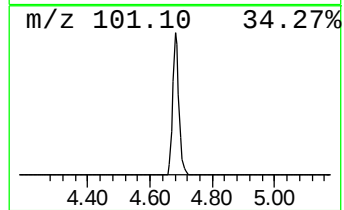
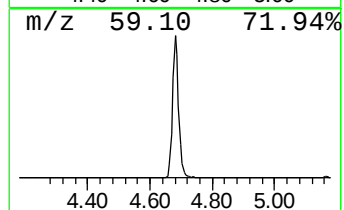
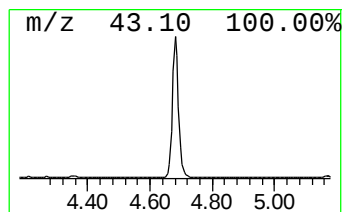
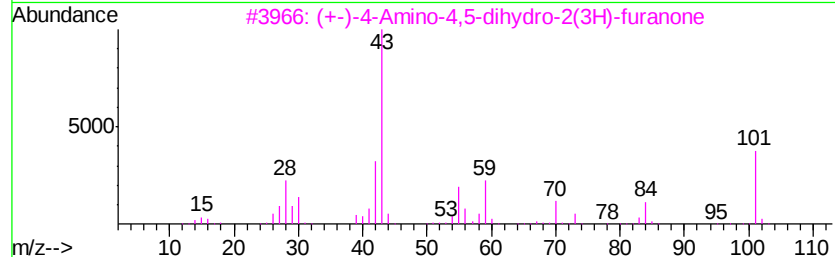
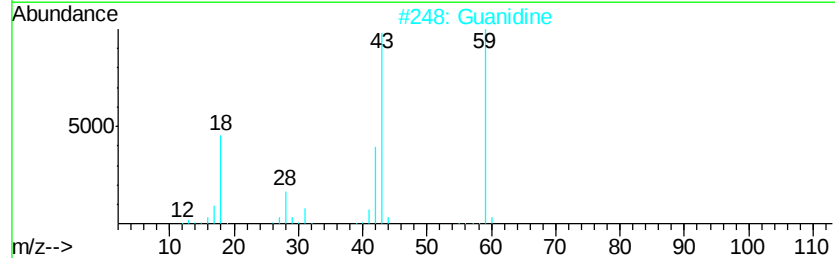
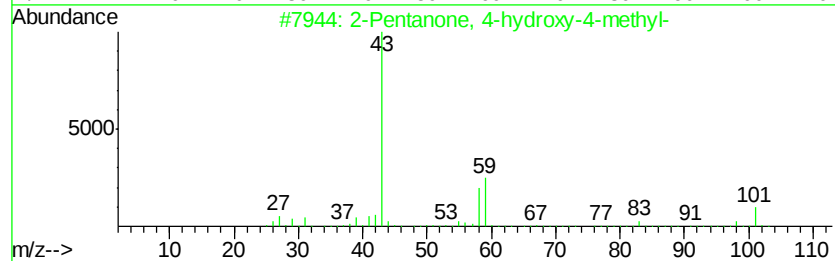
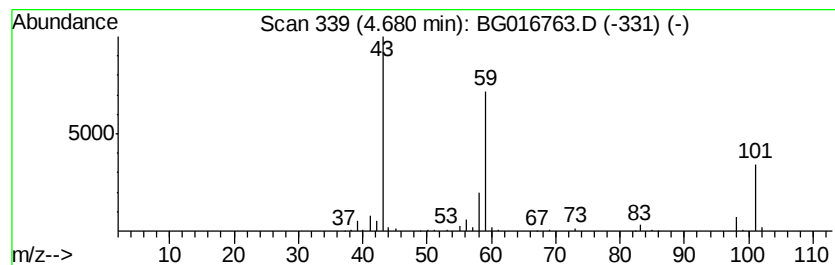
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Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.68	12.85 ng	192484	1,4-Dichlorobenzene-d4	7.57

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2		Guanidine	59	CH5N3	000113-00-8	38
3		(+)-4-Amino-4,5-dihydro-2(3H)-furanone	101	C4H7NO2	016504-58-8	38
4		1,3-Dioxolane-2-methanol, 2,4-dihydro-	132	C6H12O3	053951-43-2	33
5		4-Hydroxy-3-hexanone	116	C6H12O2	004984-85-4	25



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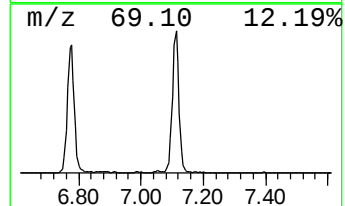
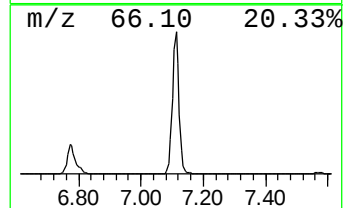
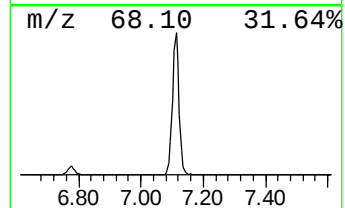
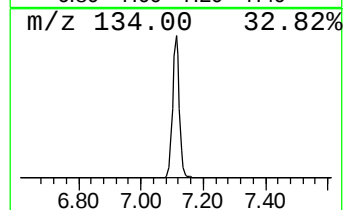
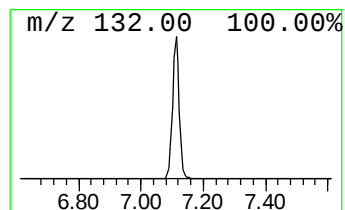
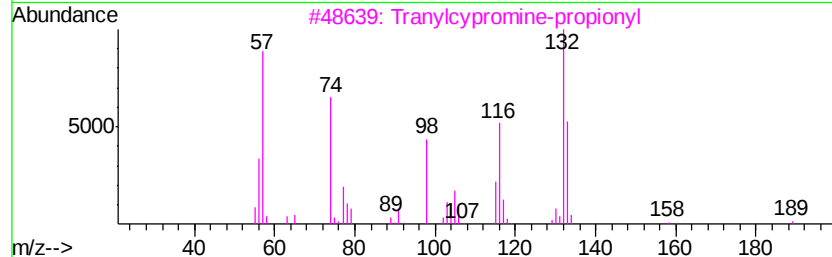
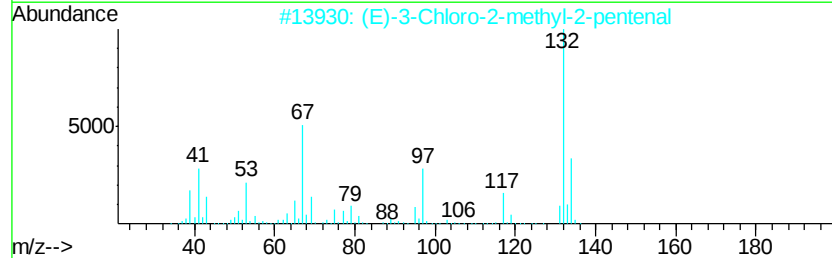
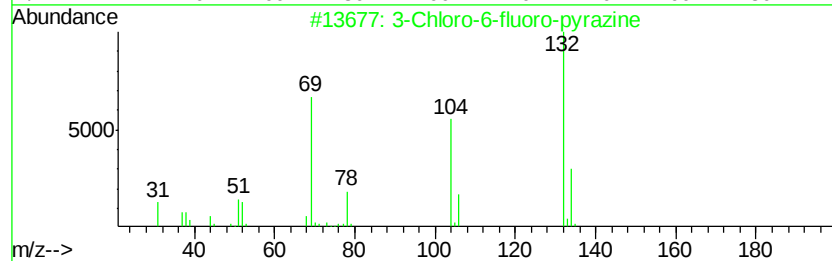
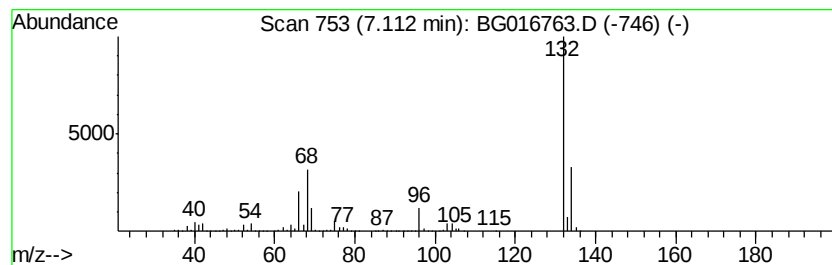
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Peak Number 3 unknown7.11 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.11	84.07 ng	1259740	1,4-Dichlorobenzene-d4	7.57

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	38
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
3		Tranvlcypromine-propionyl	189	C12H15NO	1000123-86-3	12
4		4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	9
5		4-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-60-2	9



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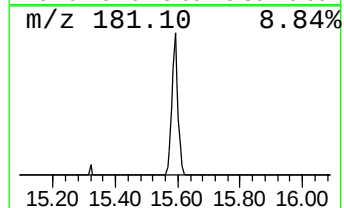
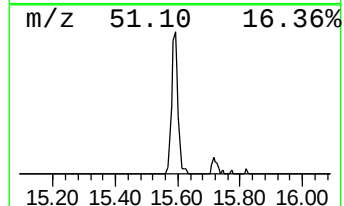
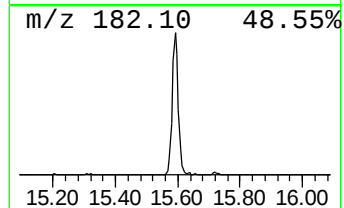
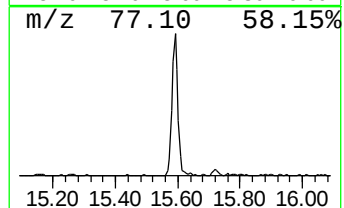
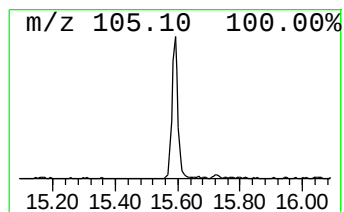
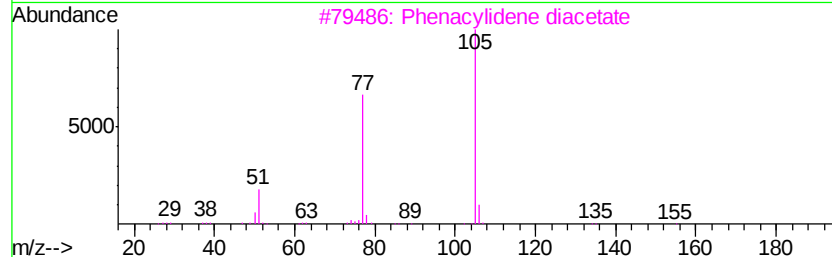
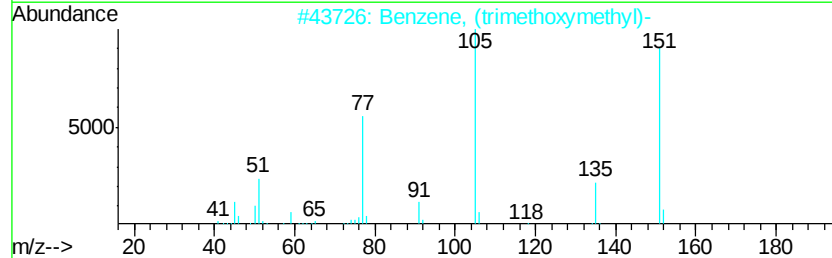
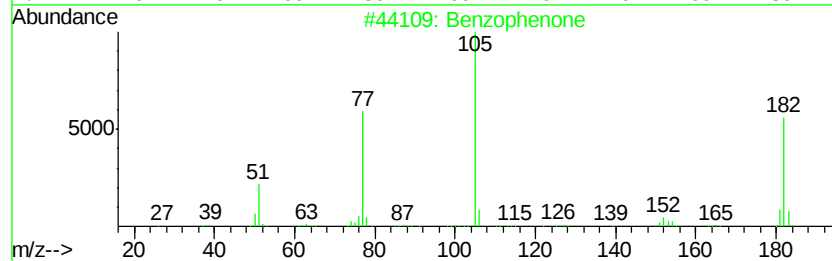
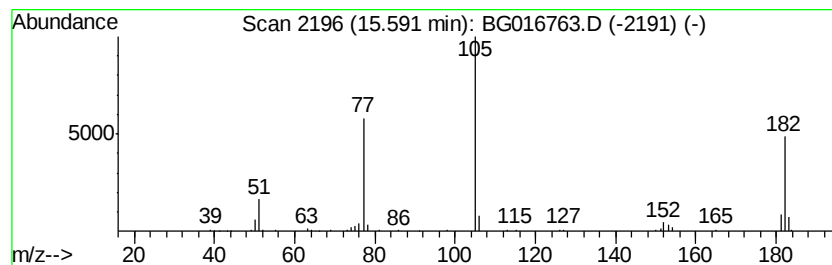
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Peak Number 5 Benzophenone Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD		R.T.
15.59	3.66 ng	114447	Acenaphthene-d10		14.23
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzophenone	182	C13H10O	000119-61-9	96
2	Benzene, (trimethoxymethyl)-	182	C10H14O3	000707-07-3	43
3	Phenacylidene diacetate	236	C12H12O5	005062-30-6	43
4	Phenylglyoxal	134	C8H6O2	001074-12-0	43
5	Benzamidomethyl benzoate	255	C15H13NO3	061652-83-3	43



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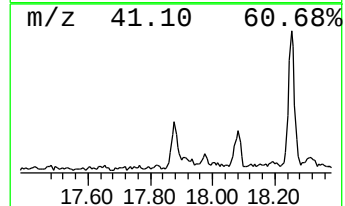
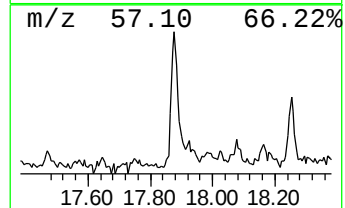
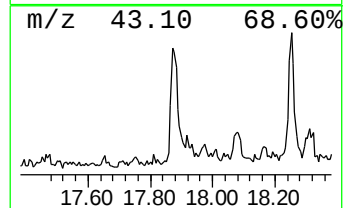
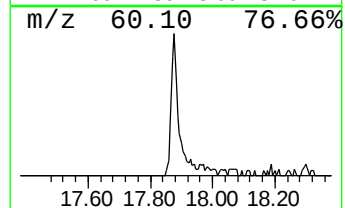
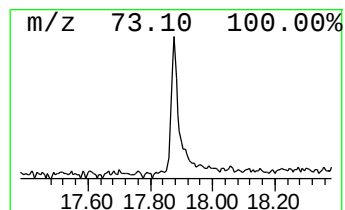
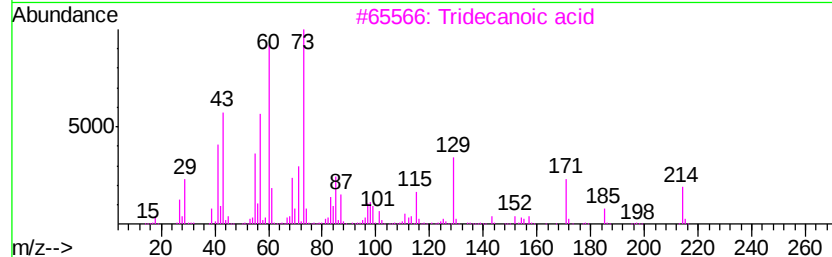
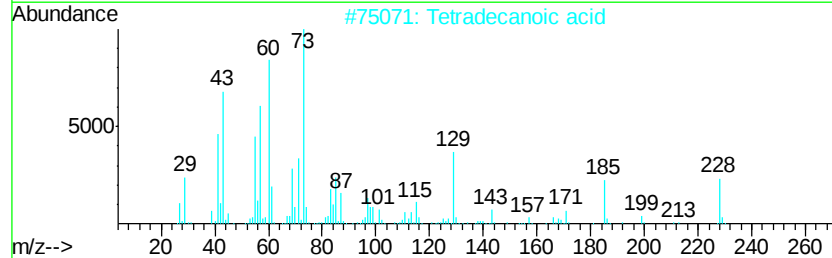
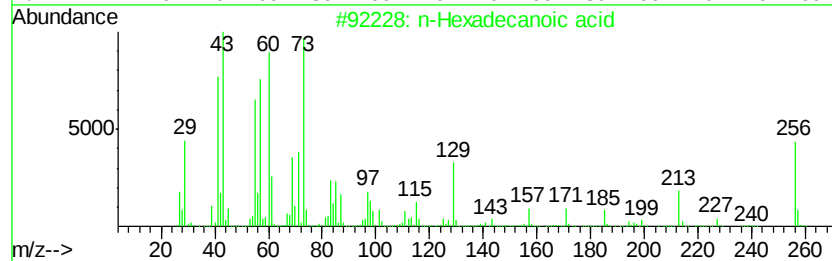
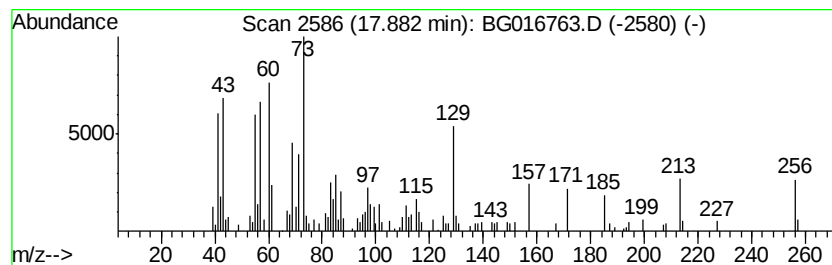
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Peak Number 6 n-Hexadecanoic acid Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.	
17.88	4.09 ng	140908	Phenanthrene-d10	16.97	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2	Tetradecanoic acid	228	C14H28O2	000544-63-8	94
3	Tridecanoic acid	214	C13H26O2	000638-53-9	76
4	Pentadecanoic acid	242	C15H30O2	001002-84-2	64
5	Undecanoic acid	186	C11H22O2	000112-37-8	50



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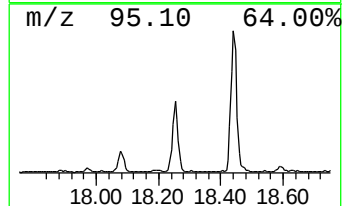
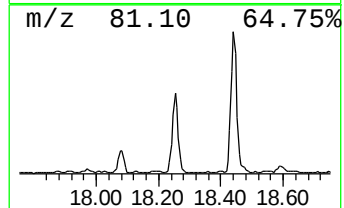
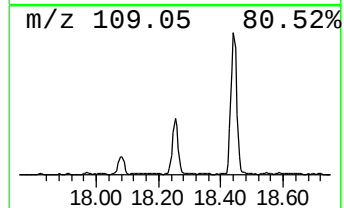
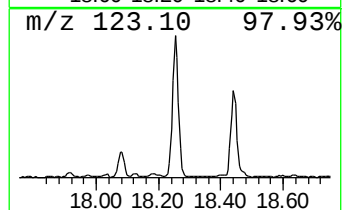
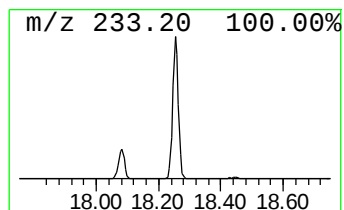
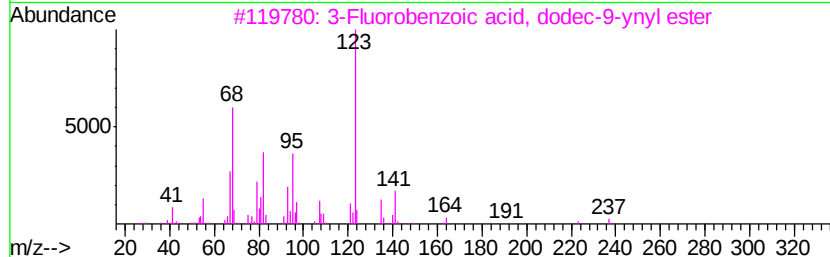
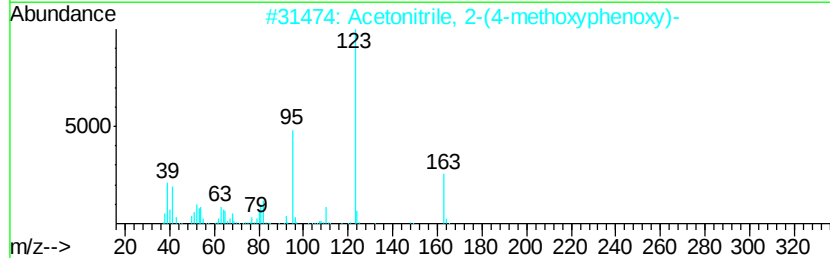
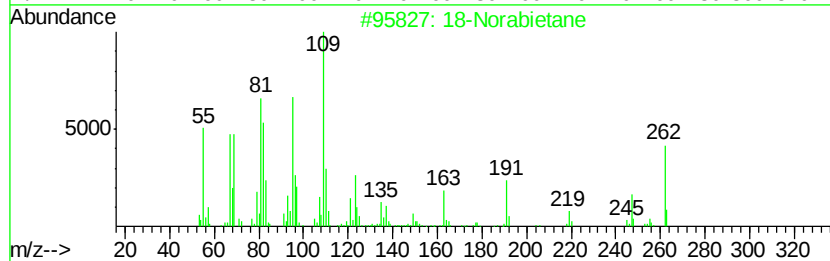
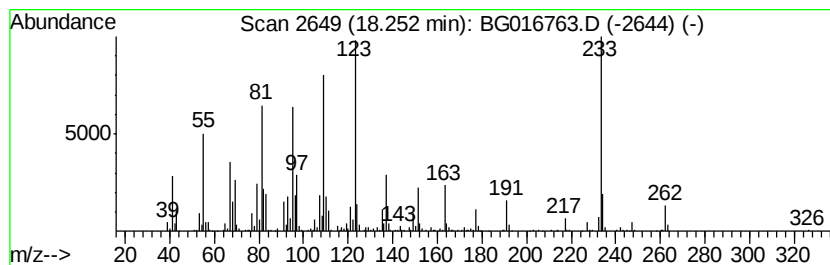
Instrument :
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ClientSampleId :
SB-74D

Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG042315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 unknown18.25 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.25	15.70 ng	540649	Phenanthrene-d10	16.97	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	18-Norabietane	262	C19H34	1000293-16-6	35
2	Acetonitrile, 2-(4-methoxyphenoxy)-	163	C9H9NO2	022446-12-4	22
3	3-Fluorobenzoic acid, dodec-9-yn...	304	C19H25F02	1000283-01-6	18
4	Bicyclo[3.1.1]hept-3-en-2-ol, 4,...	152	C10H16O	000473-67-6	14
5	3-Fluorobenzoic acid, undec-10-e...	292	C18H25F02	1000279-01-2	14



Data Path : Z:\HPCHEM1\BNA G\DATA\BG042715\
Data File : BG016763.D
Acq On : 28 Apr 2015 10:56
Operator : TP/IZ
Sample : G1991-11
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SB-74D

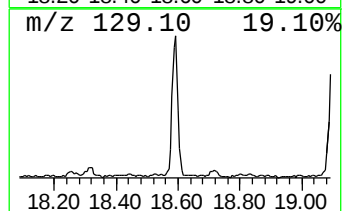
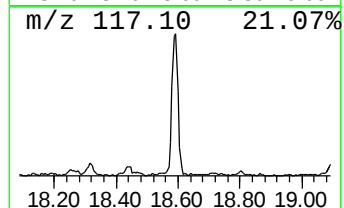
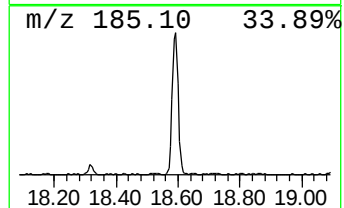
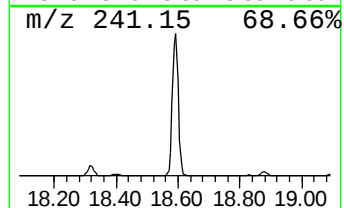
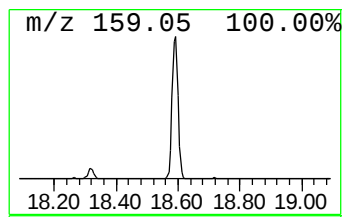
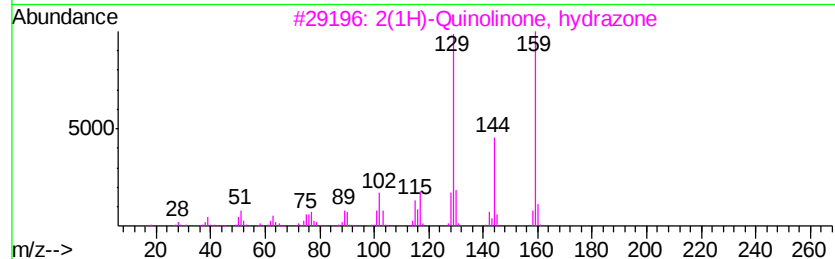
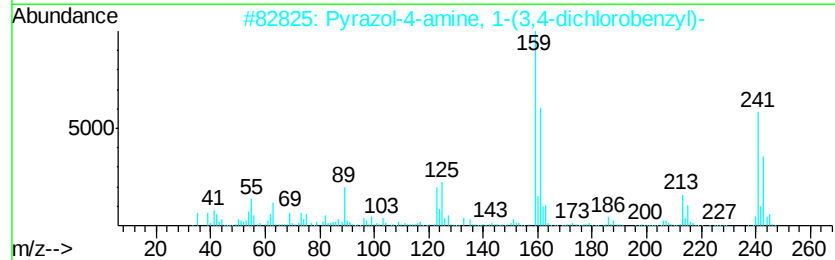
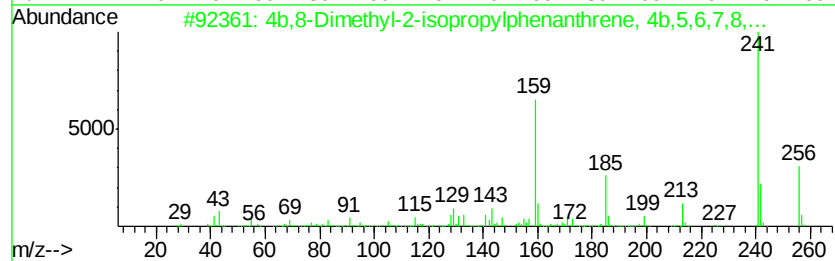
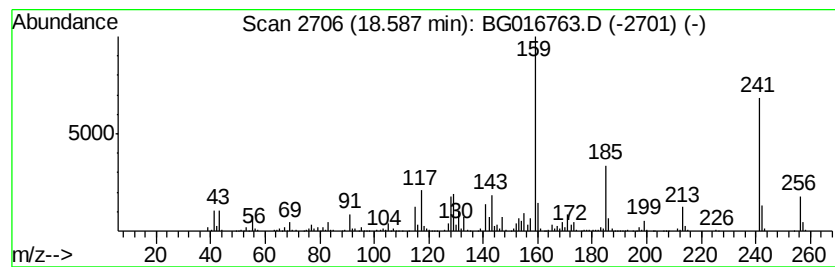
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 4b,8-Dimethyl-2-isopropylph... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.59	20.67 ng	711762	Phenanthrene-d10	16.97

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4b,8-Dimethyl-2-isopropylphenant...	256	C19H28	1000197-14-1	98
2		Pyrazol-4-amine, 1-(3,4-dichloro...	241	C10H9Cl2N3	1000273-77-4	52
3		2(1H)-Quinolinone, hvdrazone	159	C9H9N3	015793-77-8	35
4		Naphthalene, 1,2,3,4-tetrahydro-...	174	C13H18	030316-36-0	35
5		s-Indacen-1(2H)-one, 3,5,6,7-tet...	256	C18H24O	038754-94-8	30



Data Path : Z:\HPCHEM1\BNA G\DATA\BG042715\
Data File : BG016763.D
Acq On : 28 Apr 2015 10:56
Operator : TP/IZ
Sample : G1991-11
Misc :
ALS Vial : 32 Sample Multiplier: 1

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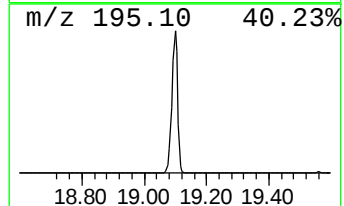
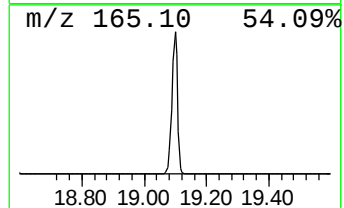
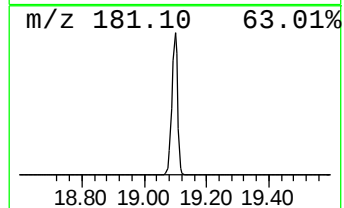
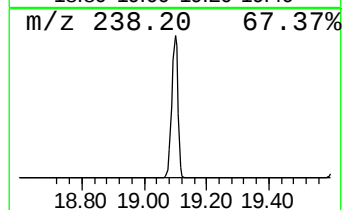
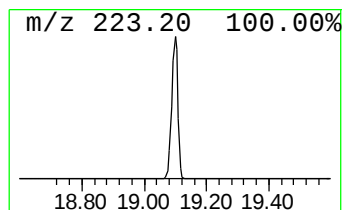
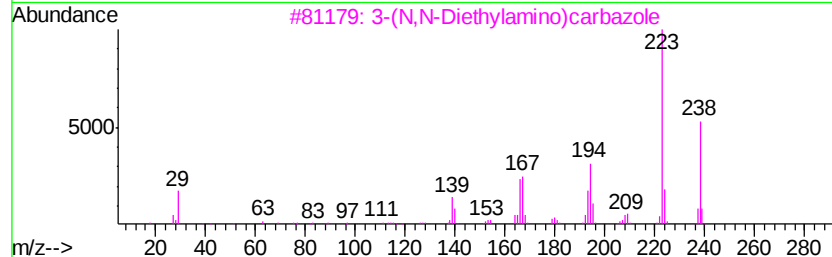
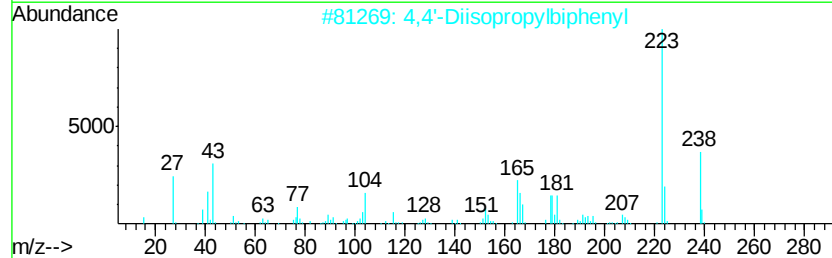
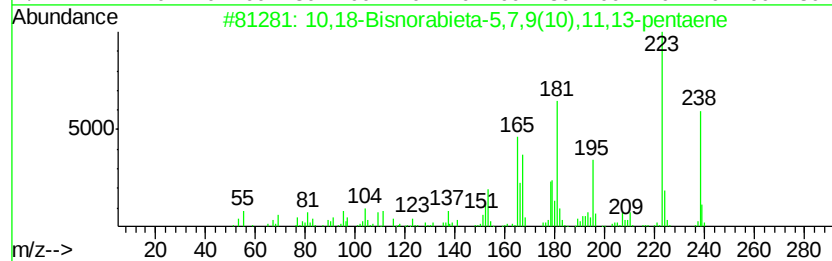
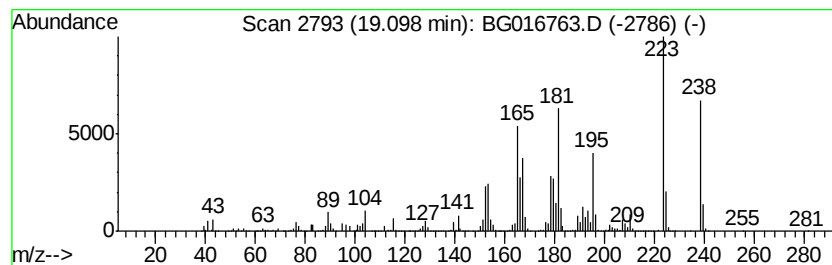
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 10,18-Bisnorabieta-5,7,9(10... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.10	224.38 ng	8295920	Chrysene-d12	21.16

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	10,18-Bisnorabieta-5,7,9(10),11,...	238	C18H22	006566-19-4	99
2		4,4'-Diisopropylbiphenyl	238	C18H22	018970-30-4	50
3		3-(N,N-Diethylamino)carbazole	238	C16H18N2	054994-31-9	49
4		9,10-Anthracenedione, 1-amino-	223	C14H9NO2	000082-45-1	45
5		3,4'-Diisopropylbiphenyl	238	C18H22	061434-46-6	43



Data Path : Z:\HPCHEM1\BNA G\DATA\BG042715\
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Acq On : 28 Apr 2015 10:56
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Sample : G1991-11
Misc :
ALS Vial : 32 Sample Multiplier: 1

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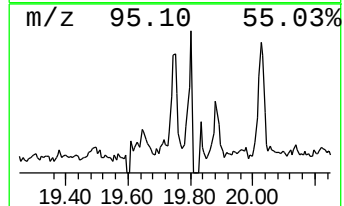
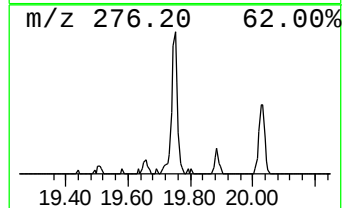
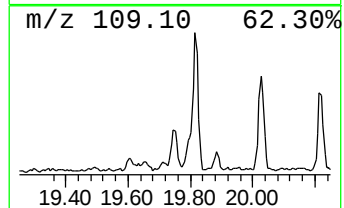
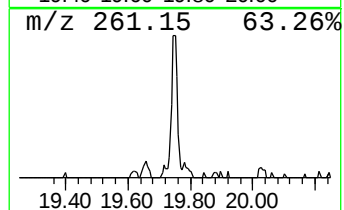
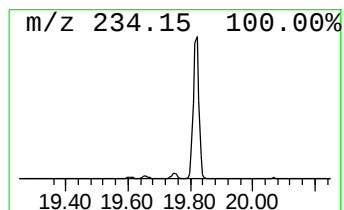
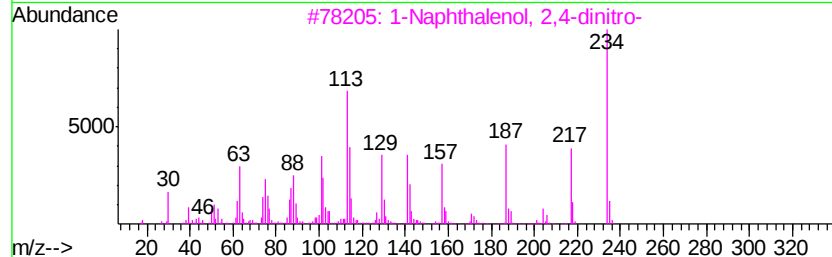
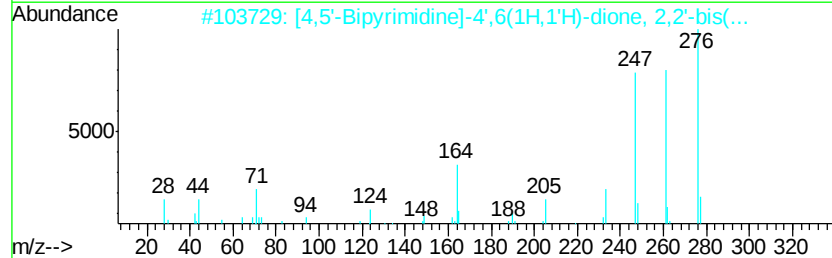
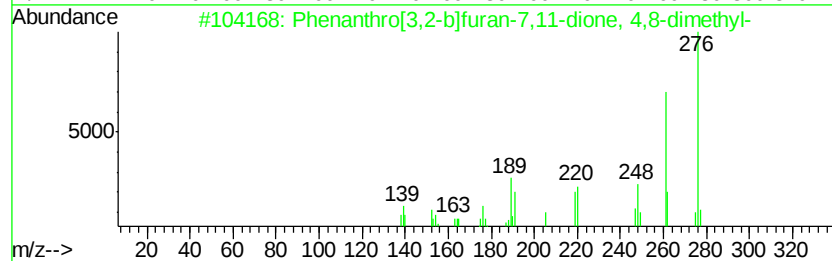
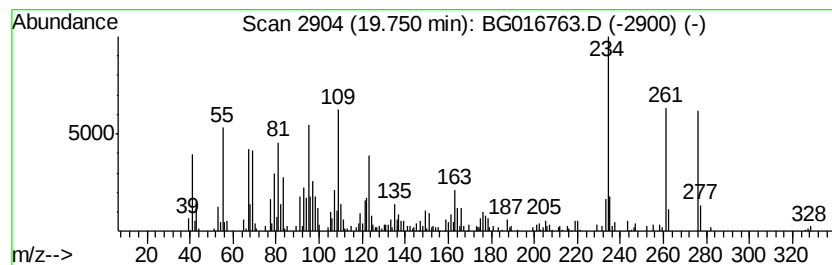
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 unknown19.75 Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.75	3.71 ng	137076	Chrysene-d12	21.16

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phenanthro[3,2-b]furan-7,11-dion...	276	C18H12O3	020958-17-2	38
2		[4,5'-Bipyrimidine]-4',6(1H,1'H)...	276	C12H16N6O2	059549-59-6	30
3		1-Naphthalenol, 2,4-dinitro-	234	C10H6N2O5	000605-69-6	25
4		6H-Pyrido[4,3-b]carbazole, 9-met...	276	C18H16N2O	010371-86-5	22
5		(1H)Pyrrole, 3,4-dimethyl-5-meth...	276	C15H20N2OS	1000197-22-6	18



Data Path : Z:\HPCHEM1\BNA G\DATA\BG042715\
Data File : BG016763.D
Acq On : 28 Apr 2015 10:56
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Sample : G1991-11
Misc :
ALS Vial : 32 Sample Multiplier: 1

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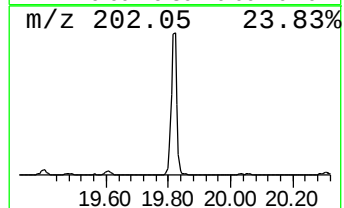
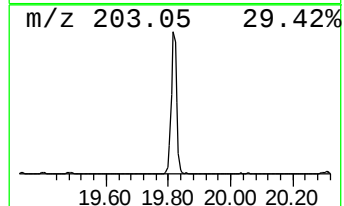
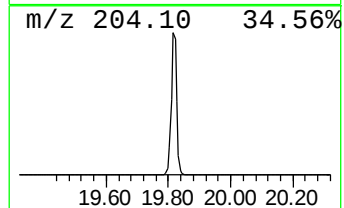
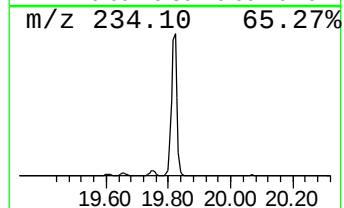
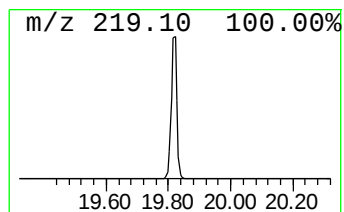
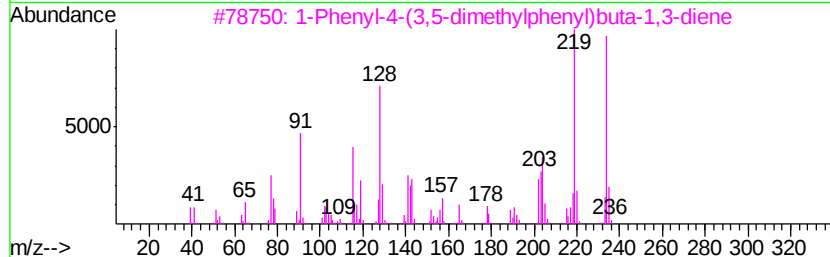
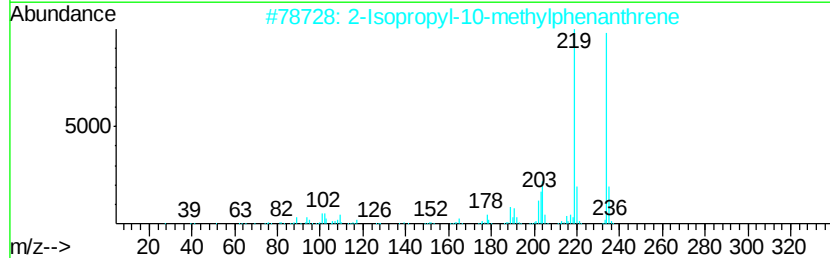
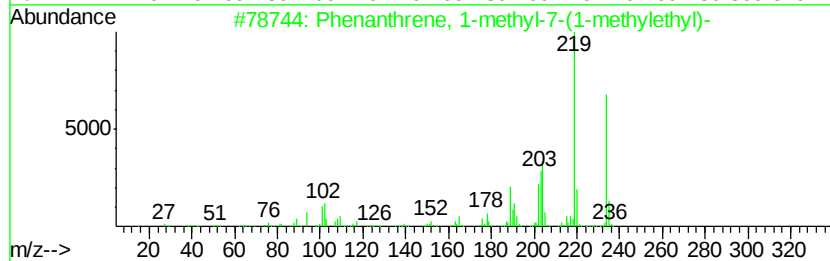
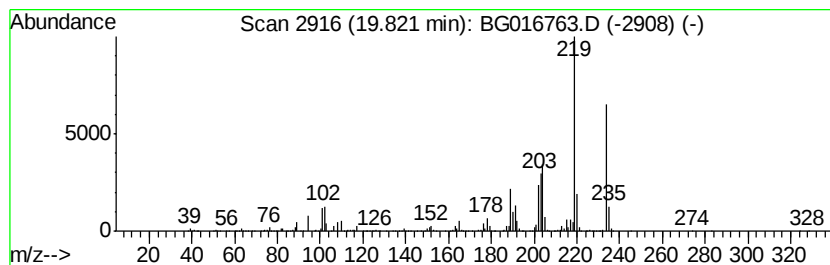
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Phenanthrene, 1-methyl-7-(1... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.82	54.13 ng	2001390	Chrysene-d12	21.16

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phenanthrene, 1-methyl-7-(1-meth...	234	C18H18	000483-65-8	99
2		2-Isopropyl-10-methylphenanthrene	234	C18H18	066552-97-4	94
3		1-Phenyl-4-(3,5-dimethylphenyl)b...	234	C18H18	1000202-15-0	80
4		Phenanthrene, 2,4,5,7-tetramethyl-	234	C18H18	007396-38-5	70
5		1-(10-Methylantracen-9-yl)ethanone	234	C17H14O	036778-18-4	70



Data Path : Z:\HPCHEM1\BNA G\DATA\BG042715\
Data File : BG016763.D
Acq On : 28 Apr 2015 10:56
Operator : TP/IZ
Sample : G1991-11
Misc :
ALS Vial : 32 Sample Multiplier: 1

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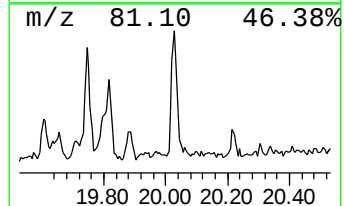
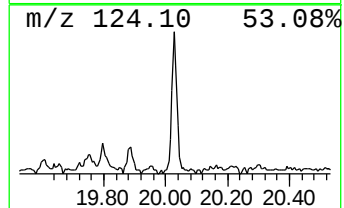
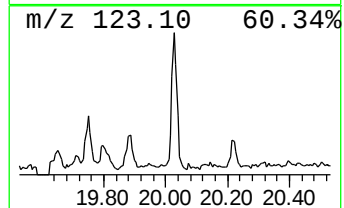
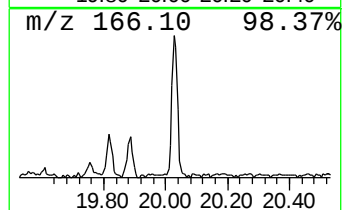
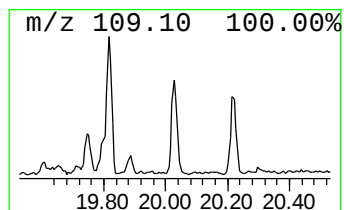
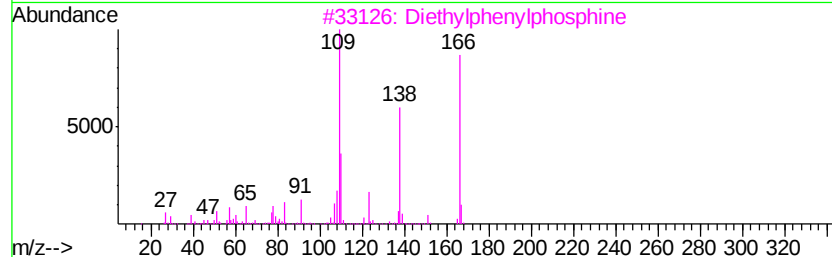
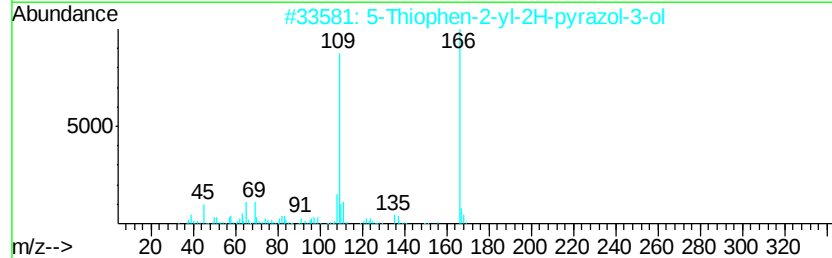
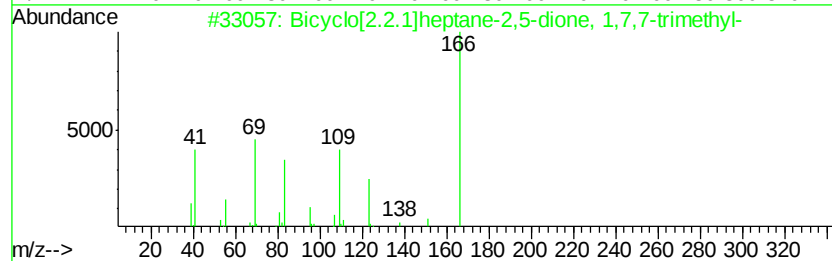
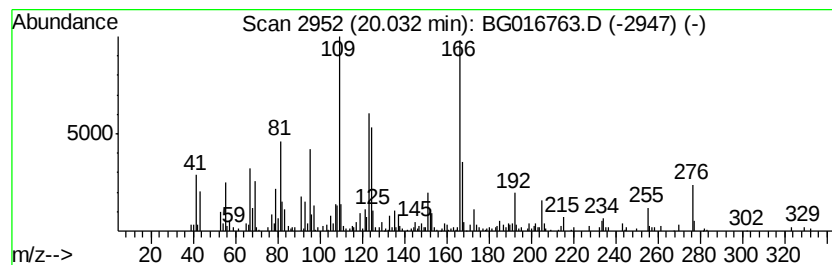
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 unknown20.03 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.03	4.65 ng	171741	Chrysene-d12	21.16

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Bicyclo[2.2.1]heptane-2,5-dione,...	166	C10H14O2	004230-32-4	46
2		5-Thiophen-2-yl-2H-pyrazol-3-ol	166	C7H6N2OS	1000278-27-0	35
3		Diethylphenylphosphine	166	C10H15P	001605-53-4	35
4		Ethanone, 1-(2-methyl-1-cyclopent...	124	C8H12O	003168-90-9	18
5		Durohydroquinone	166	C10H14O2	000527-18-4	18



Data Path : Z:\HPCHEM1\BNA G\DATA\BG042715\
Data File : BG016763.D
Acq On : 28 Apr 2015 10:56
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Misc :
ALS Vial : 32 Sample Multiplier: 1

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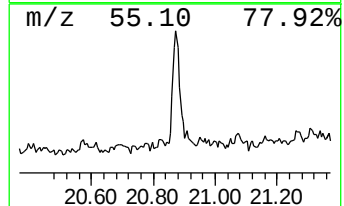
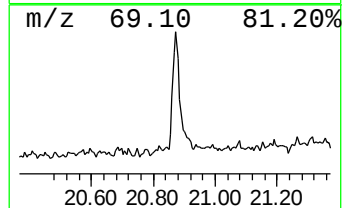
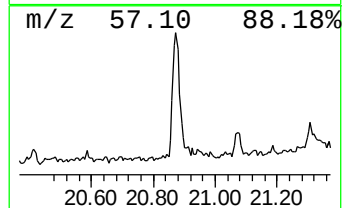
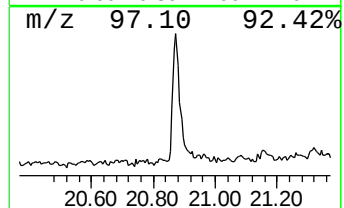
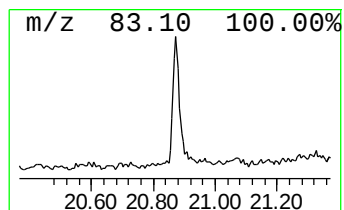
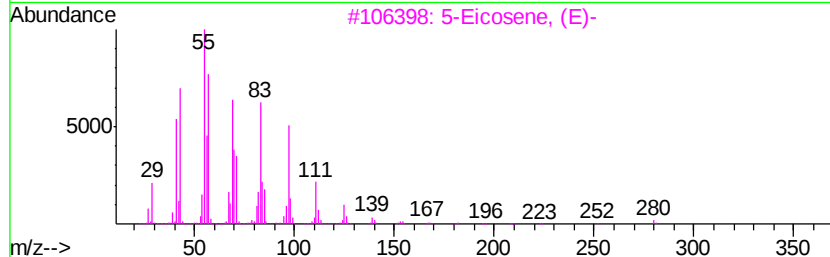
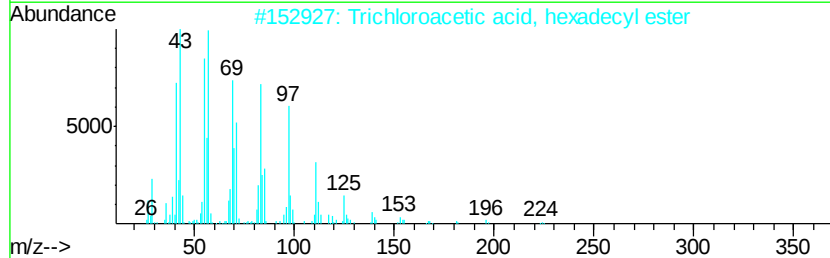
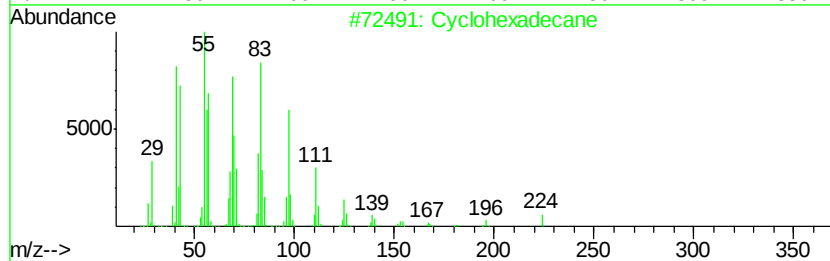
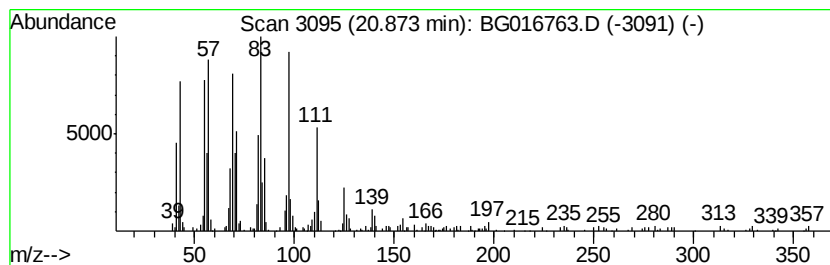
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Cyclohexadecane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.87	4.77 ng	176189	Chrysene-d12	21.16

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexadecane	224	C16H32	000295-65-8	93
2		Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	90
3		5-Eicosene, (E)-	280	C20H40	074685-30-6	90
4		1-Hexacosanol	382	C26H54O	000506-52-5	83
5		1-Docosene	308	C22H44	001599-67-3	76



Data Path : Z:\HPCHEM1\BNA G\DATA\BG042715\
Data File : BG016763.D
Acq On : 28 Apr 2015 10:56
Operator : TP/IZ
Sample : G1991-11
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SB-74D

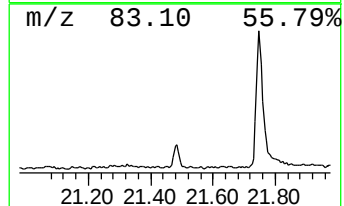
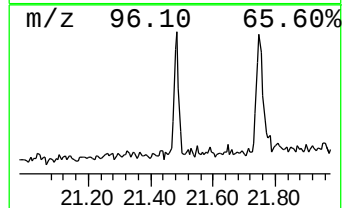
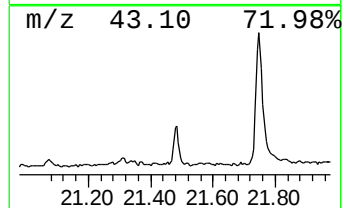
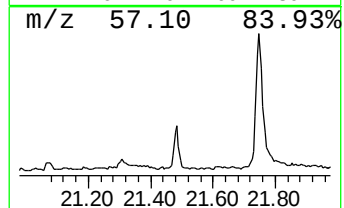
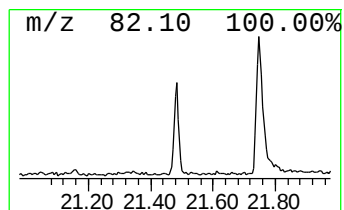
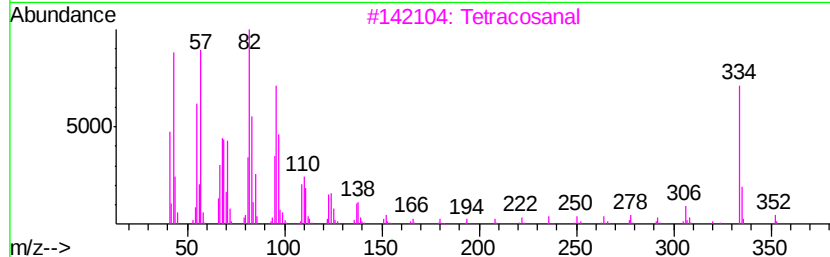
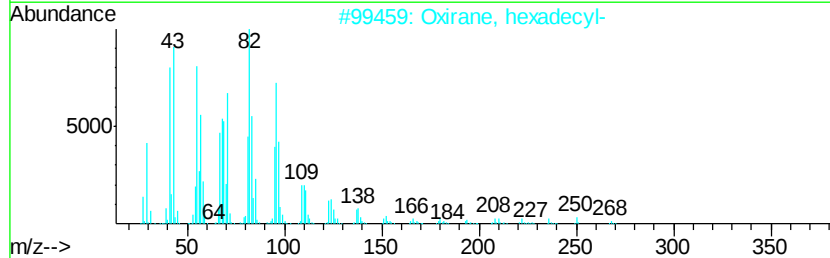
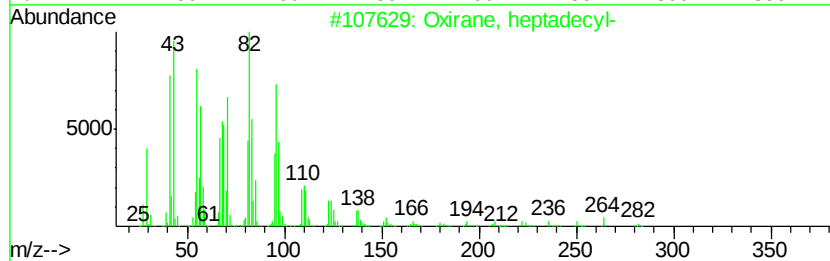
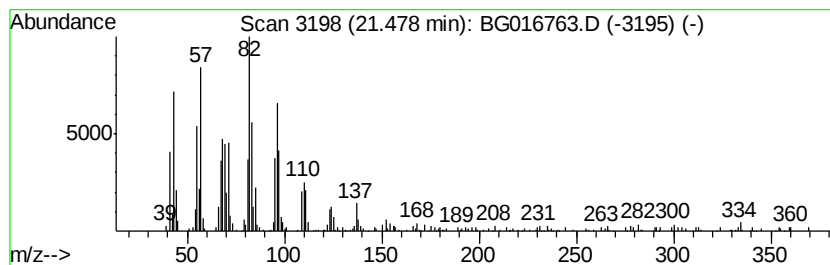
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 Oxirane, heptadecyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.48	3.93 ng	145121	Chrysene-d12	21.16

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Oxirane, heptadecyl-	282	C19H38O	067860-04-2	96
2		Oxirane, hexadecyl-	268	C18H36O	007390-81-0	91
3		Tetracosanal	352	C24H48O	057866-08-7	91
4		1,21-Docosadiene	306	C22H42	053057-53-7	91
5		Hexadecanal	240	C16H32O	000629-80-1	91



Data Path : Z:\HPCHEM1\BNA G\DATA\BG042715\
Data File : BG016763.D
Acq On : 28 Apr 2015 10:56
Operator : TP/IZ
Sample : G1991-11
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SB-74D

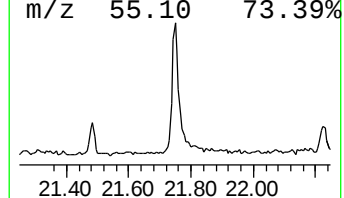
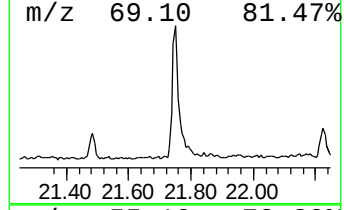
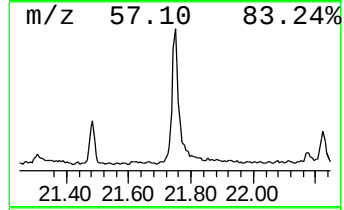
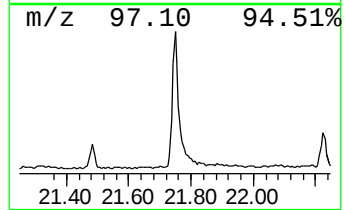
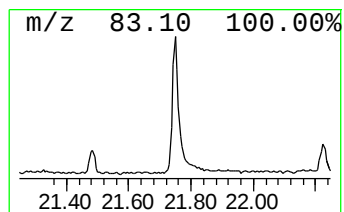
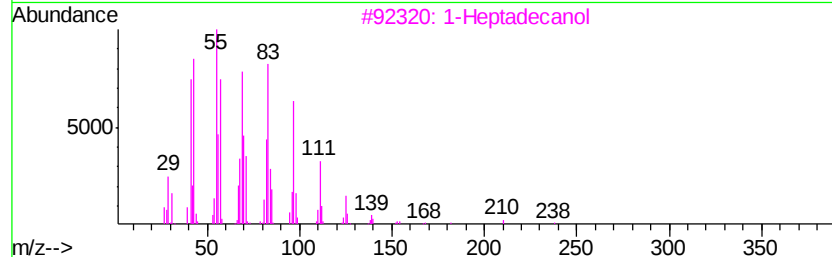
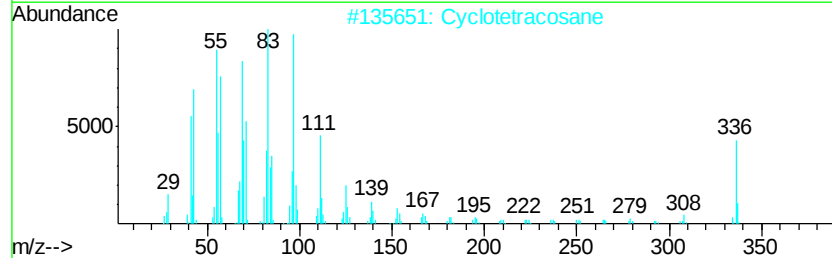
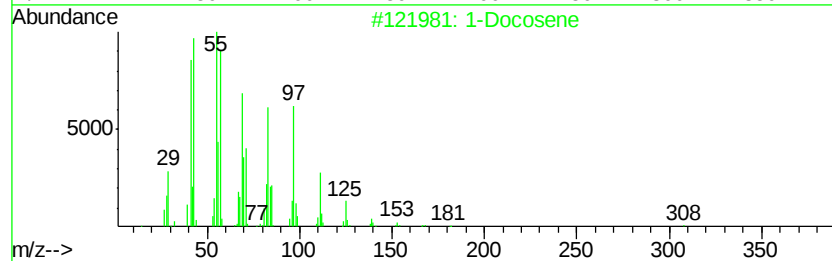
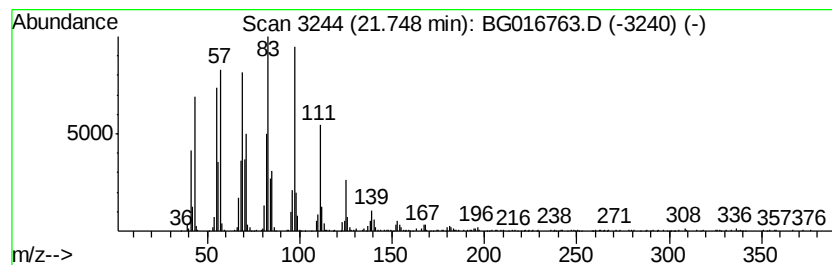
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 1-Docosene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.75	15.94 ng	589364	Chrysene-d12	21.16

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Docosene	308	C22H44	001599-67-3	93
2		Cyclotetracosane	336	C24H48	000297-03-0	93
3		1-Heptadecanol	256	C17H36O	001454-85-9	91
4		1-Heneicosyl formate	340	C22H44O2	077899-03-7	91
5		3-Eicosene, (E)-	280	C20H40	074685-33-9	90



Data Path : Z:\HPCHEM1\BNA G\DATA\BG042715\
Data File : BG016763.D
Acq On : 28 Apr 2015 10:56
Operator : TP/IZ
Sample : G1991-11
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SB-74D

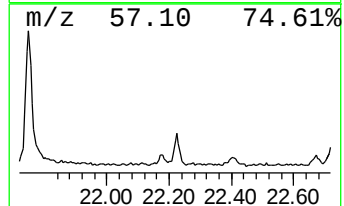
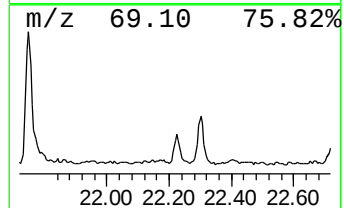
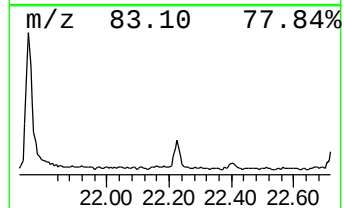
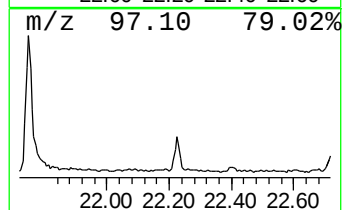
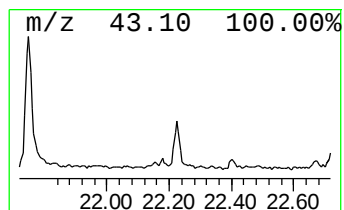
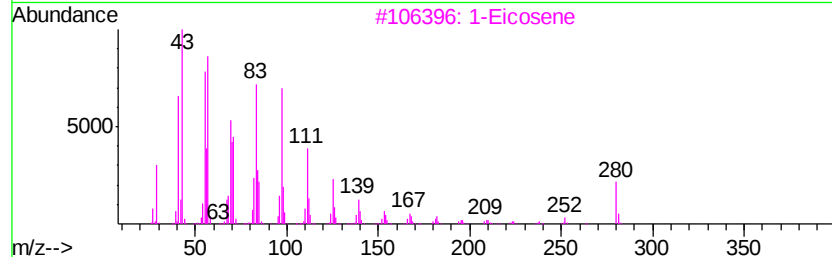
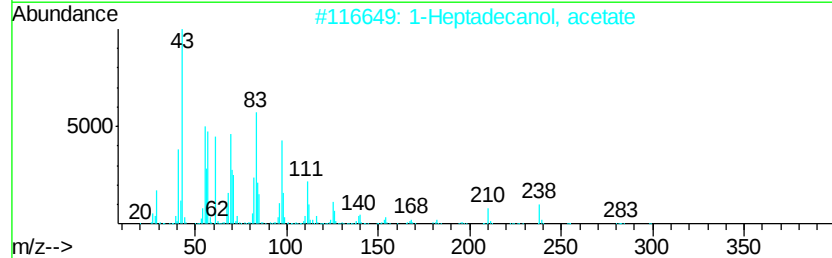
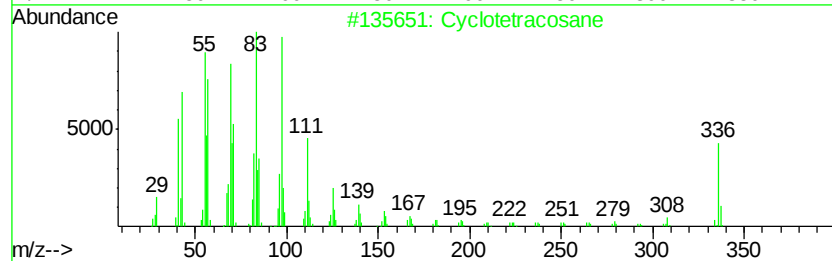
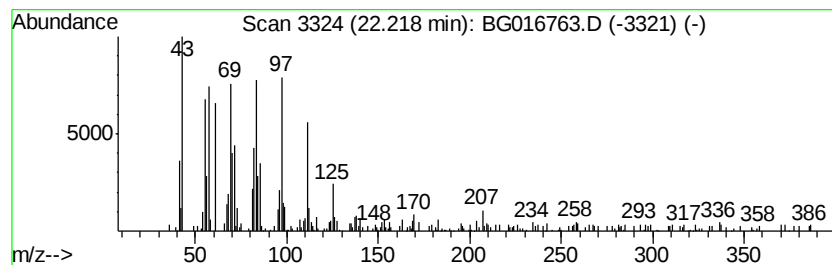
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 Cyclotetracosane Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.22	3.78 ng	139932	Chrysene-d12	21.16

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclotetracosane	336	C24H48	000297-03-0	96
2		1-Heptadecanol, acetate	298	C19H38O2	000822-20-8	93
3		1-Eicosene	280	C20H40	003452-07-1	93
4		1-Heneicosanol	312	C21H44O	015594-90-8	90
5		1-Docosene	308	C22H44	001599-67-3	87



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG042715\
Data File : BG016763.D
Acq On : 28 Apr 2015 10:56
Operator : TP/IZ
Sample : G1991-11
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SB-74D

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG042315.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
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2-Pentanone, 4-hy...	4.68	12.9	ng	192484	1	7.57	299681	20.0
unknown7.11	7.11	84.1	ng	1259740	1	7.57	299681	20.0
Benzophenone	15.59	3.7	ng	114447	3	14.23	625770	20.0
n-Hexadecanoic acid	17.88	4.1	ng	140908	4	16.97	688638	20.0
unknown18.25	18.25	15.7	ng	540649	4	16.97	688638	20.0
4b,8-Dimethyl-2-i...	18.59	20.7	ng	711762	4	16.97	688638	20.0
10,18-Bisnorabiet...	19.10	224.4	ng	8295920	5	21.16	739447	20.0
unknown19.75	19.75	3.7	ng	137076	5	21.16	739447	20.0
Phenanthrene, 1-m...	19.82	54.1	ng	2001390	5	21.16	739447	20.0
unknown20.03	20.03	4.7	ng	171741	5	21.16	739447	20.0
Cyclohexadecane	20.87	4.8	ng	176189	5	21.16	739447	20.0
Oxirane, heptadecyl-	21.48	3.9	ng	145121	5	21.16	739447	20.0
1-Docosene	21.75	15.9	ng	589364	5	21.16	739447	20.0
Cyclotetracosane	22.22	3.8	ng	139932	5	21.16	739447	20.0