

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG092316\
 Data File : BG024114.D
 Acq On : 24 Sep 2016 8:11
 Operator : UM/SJ
 Sample : H4893-05
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PZ-8

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

Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG092316.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	5.104	382	386	393	rVB2	71431	116477	2.17%	0.455%
2	5.586	463	468	485	rVB	510521	933567	17.43%	3.646%
3	7.213	739	745	758	rBV	341504	669288	12.49%	2.614%
4	7.572	799	806	820	rBV	1097238	1952311	36.44%	7.625%
5	8.036	879	885	894	rBV	195129	349033	6.52%	1.363%
6	8.365	933	941	950	rBV	942450	1641942	30.65%	6.413%
7	9.222	1079	1087	1099	rBV	888091	1674773	31.26%	6.541%
8	10.873	1361	1368	1381	rBV	283080	550766	10.28%	2.151%
9	12.124	1575	1581	1588	rVB2	36702	63206	1.18%	0.247%
10	13.229	1763	1769	1774	rBV3	41208	76838	1.43%	0.300%
11	13.329	1776	1786	1794	rBV	2426916	3762652	70.24%	14.696%
12	13.722	1849	1853	1860	rVB7	30909	53748	1.00%	0.210%
13	14.715	2014	2022	2033	rVB	516982	872151	16.28%	3.406%
14	14.809	2033	2038	2042	rBV8	26336	59239	1.11%	0.231%
15	16.095	2253	2257	2266	rVB8	34060	78003	1.46%	0.305%
16	16.219	2272	2278	2288	rBV	2245146	3458172	64.55%	13.507%
17	17.206	2441	2446	2450	rBV7	37039	69485	1.30%	0.271%
18	17.253	2452	2454	2459	rVB6	40488	57346	1.07%	0.224%
19	17.482	2487	2493	2501	rBV	650662	1001771	18.70%	3.913%
20	18.357	2638	2642	2648	rBV5	88871	143707	2.68%	0.561%
21	19.626	2854	2858	2866	rBV6	106486	166344	3.11%	0.650%
22	19.761	2877	2881	2885	rVB2	102158	120821	2.26%	0.472%
23	20.090	2931	2937	2944	rBV	4146248	5357022	100.00%	20.923%
24	20.807	3055	3059	3061	rBV	69071	75560	1.41%	0.295%
25	20.977	3085	3088	3093	rVB	91360	104045	1.94%	0.406%
26	21.788	3220	3226	3237	rVB	656311	1141239	21.30%	4.457%
27	25.124	3785	3794	3805	rVB2	347242	1054094	19.68%	4.117%

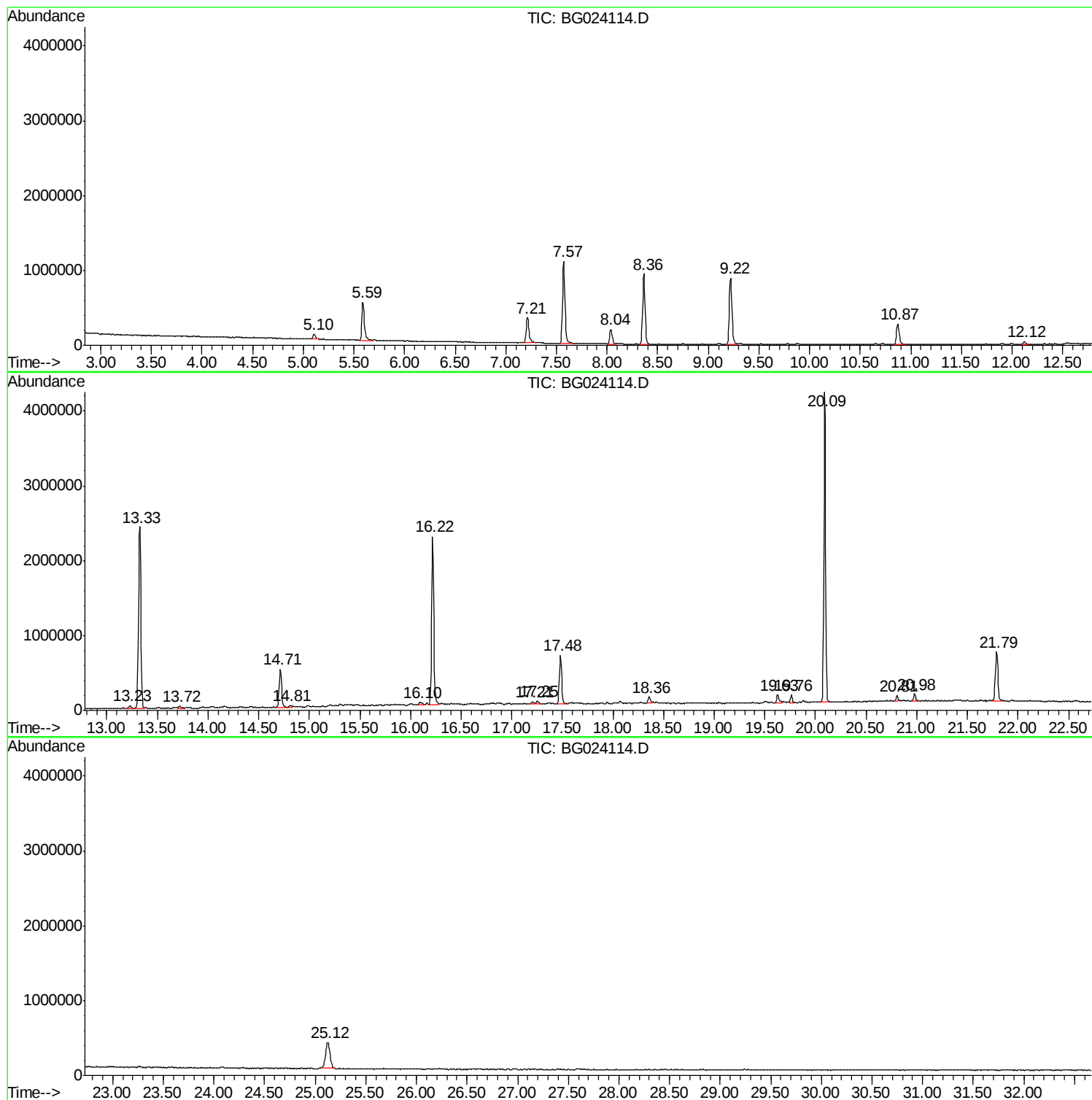
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Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG092316.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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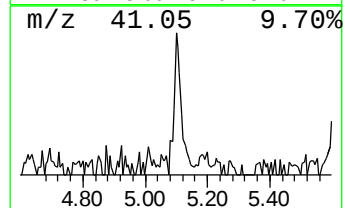
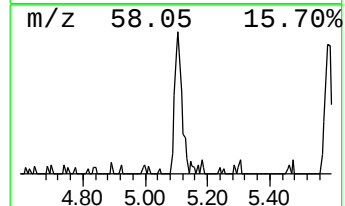
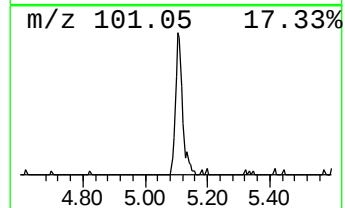
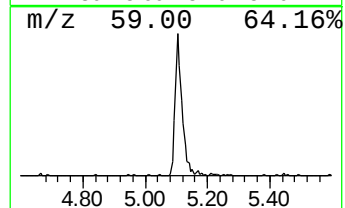
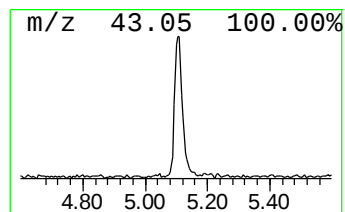
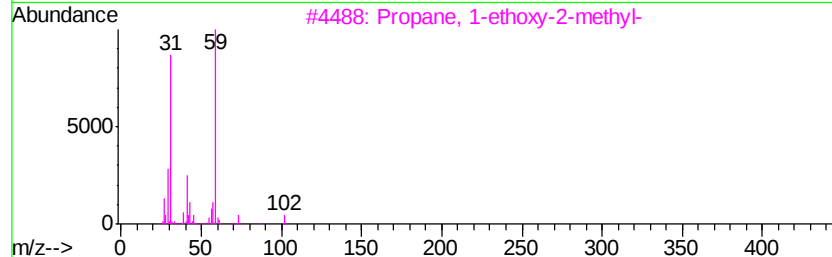
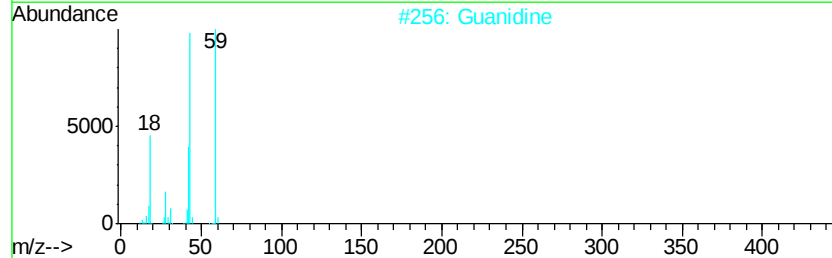
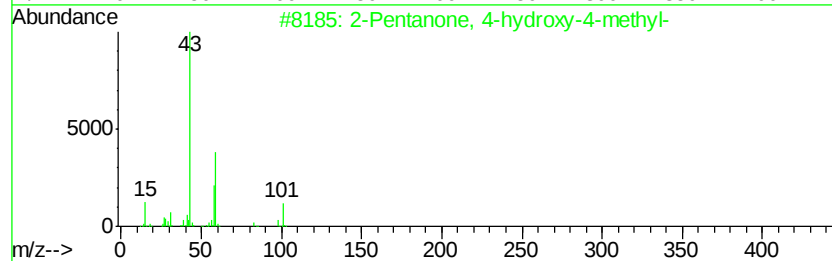
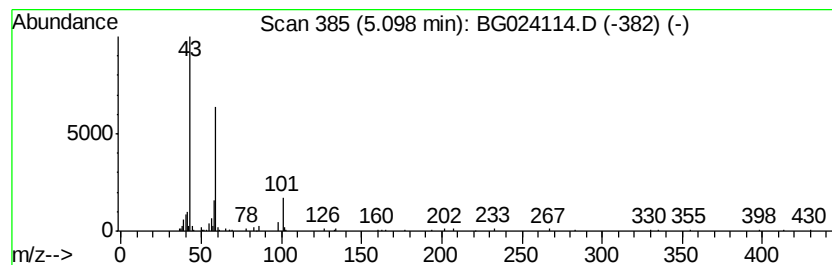
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.10	6.67 ng	116477	1,4-Dichlorobenzene-d4	8.04

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2			Guanidine	59	CH5N3	000113-00-8	50
3			Propane, 1-ethoxy-2-methyl-	102	C6H14O	000627-02-1	43
4			Dimethadione	129	C5H7NO3	000695-53-4	38
5			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	37



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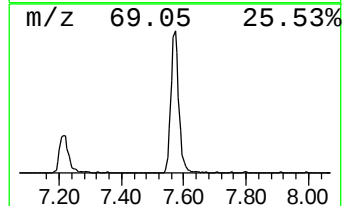
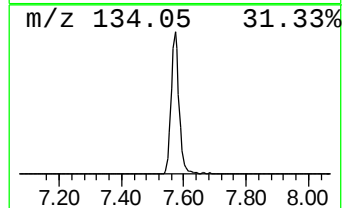
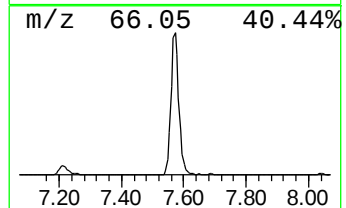
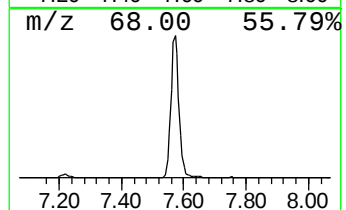
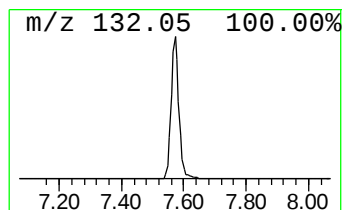
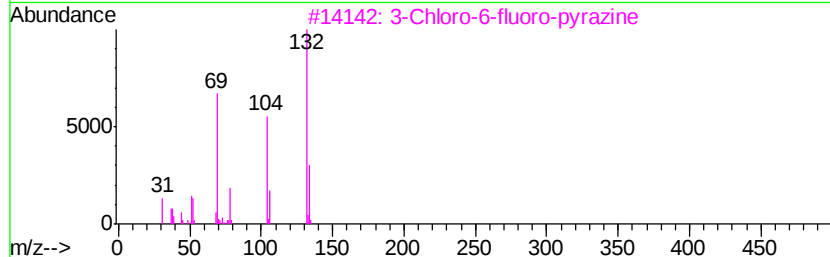
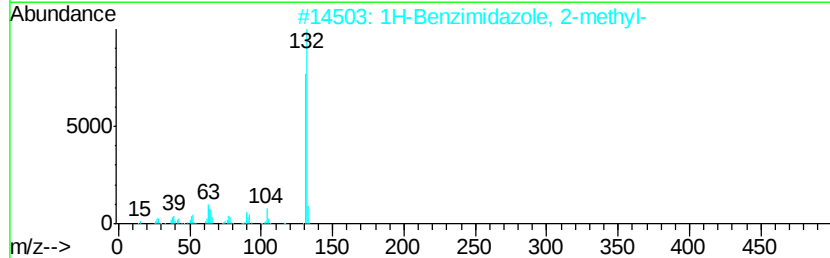
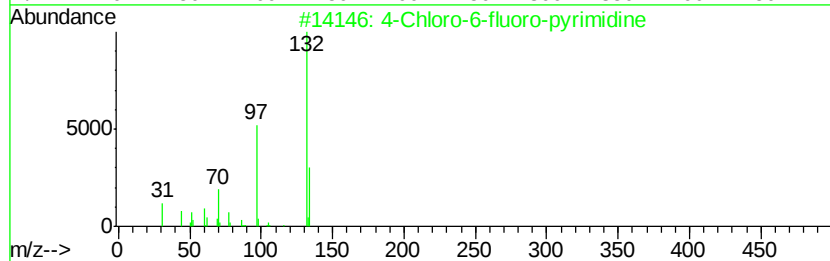
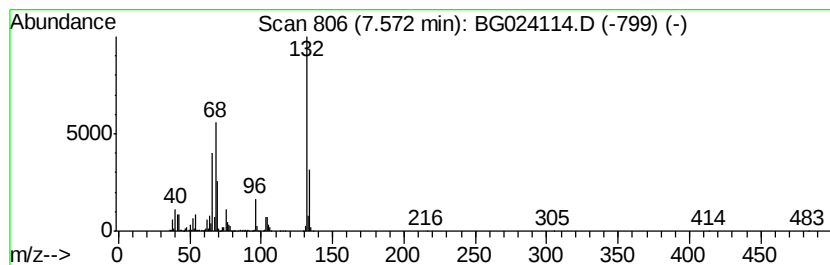
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 2 unknown7.57 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.57	111.87 ng	1952310	1,4-Dichlorobenzene-d4	8.04

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	14
2		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	14
3		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	14
4		Imidazole, 4,5-dicyano-1-methyl-	132	C6H4N4	019485-35-9	11
5		1-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-59-9	9



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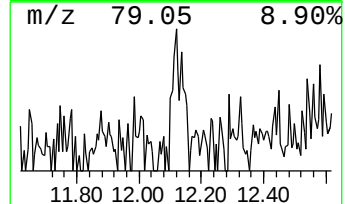
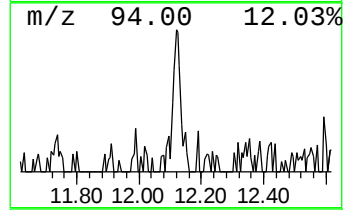
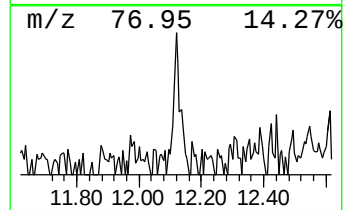
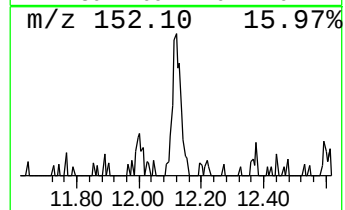
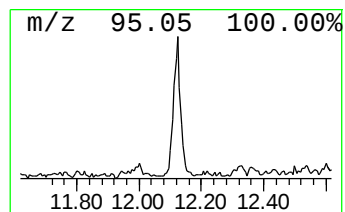
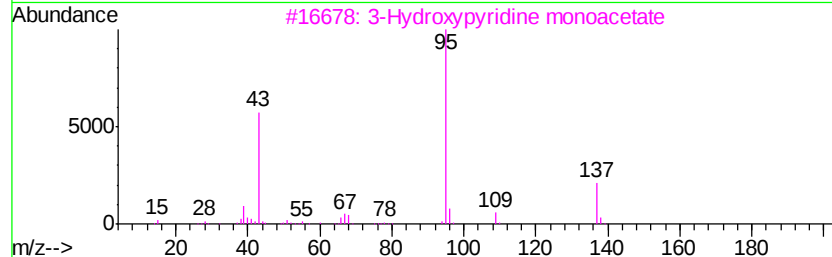
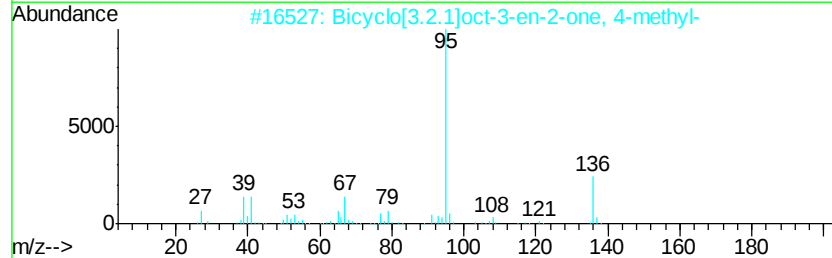
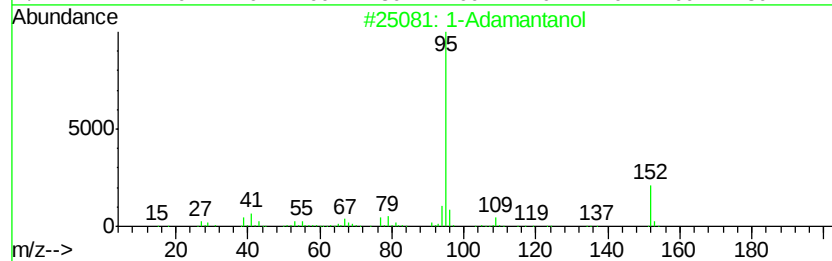
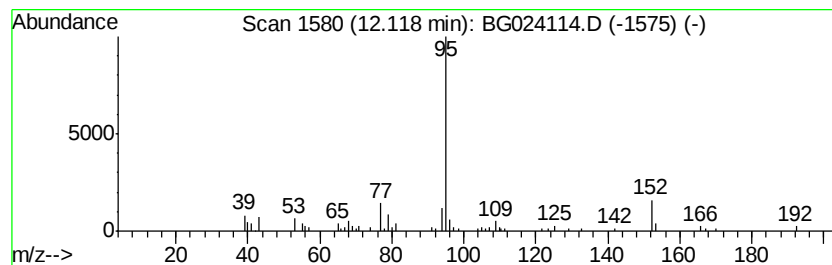
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Peak Number 4 1-Adamantanol Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.12	2.30 ng	63206	Naphthalene-d8	10.87

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Adamantanol		152	C10H16O	000768-95-6	78
2	Bicyclo[3.2.1]oct-3-en-2-one, 4-...		136	C9H12O	062702-89-0	53
3	3-Hydroxypyridine monoacetate		137	C7H7NO2	017747-43-2	53
4	1,3-Dimethyl-1-cyclohexene		110	C8H14	002808-76-6	50
5	2-Pyrimidinamine		95	C4H5N3	000109-12-6	42



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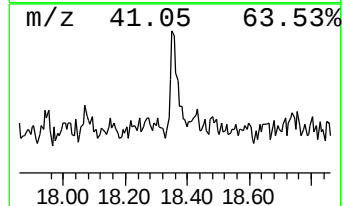
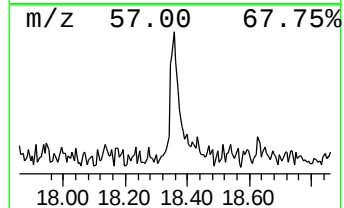
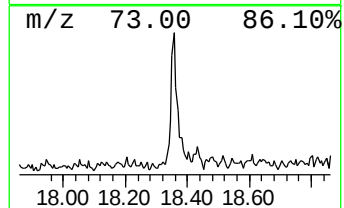
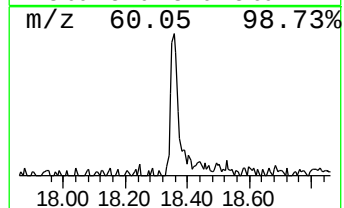
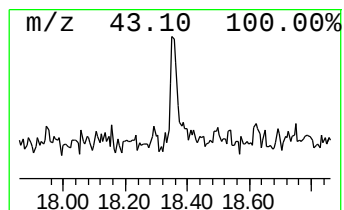
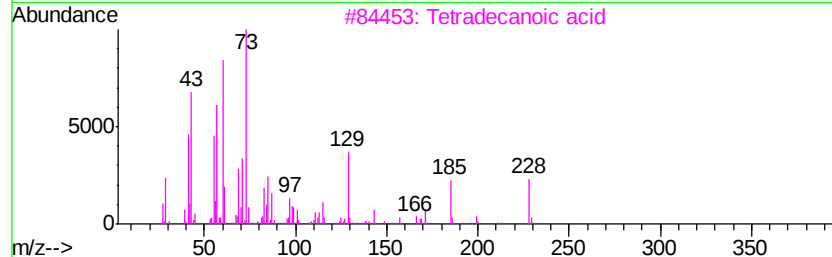
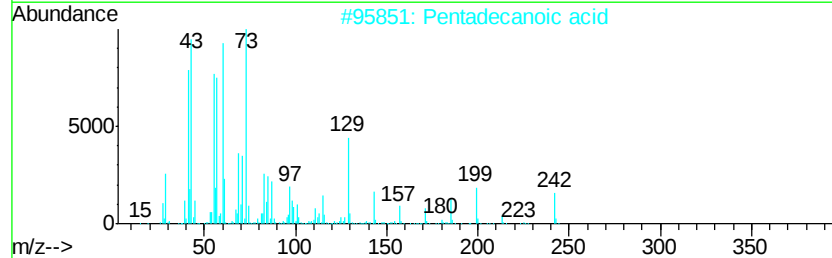
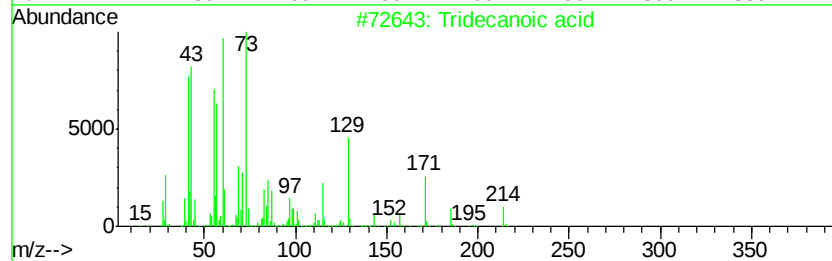
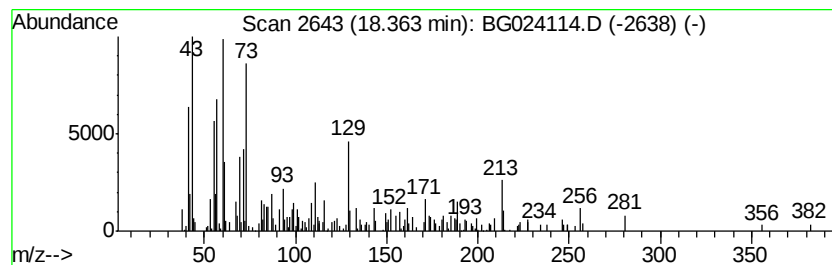
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 5 Tridecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.36	2.87 ng	143707	Phenanthrene-d10	17.48

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tridecanoic acid	214	C13H26O2	000638-53-9	81
2		Pentadecanoic acid	242	C15H30O2	001002-84-2	74
3		Tetradecanoic acid	228	C14H28O2	000544-63-8	68
4		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	59
5		Octadecanoic acid	284	C18H36O2	000057-11-4	59



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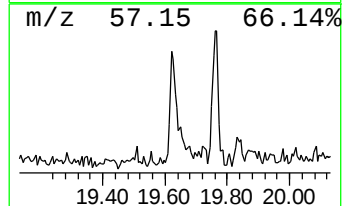
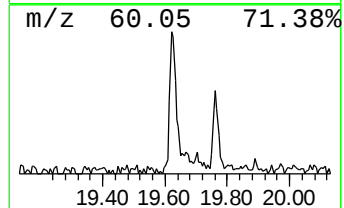
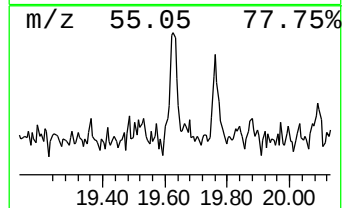
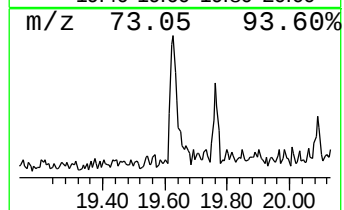
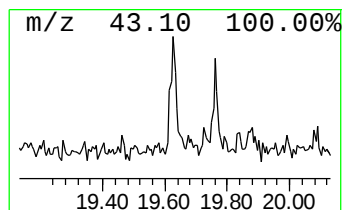
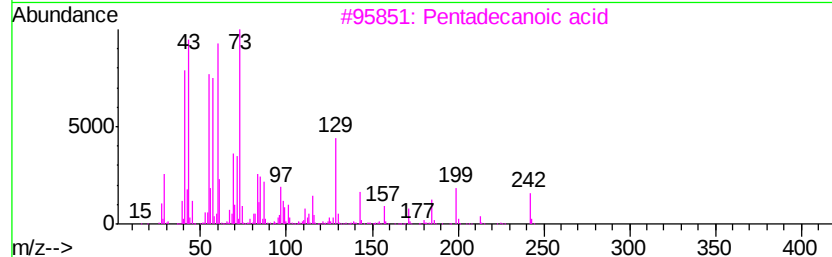
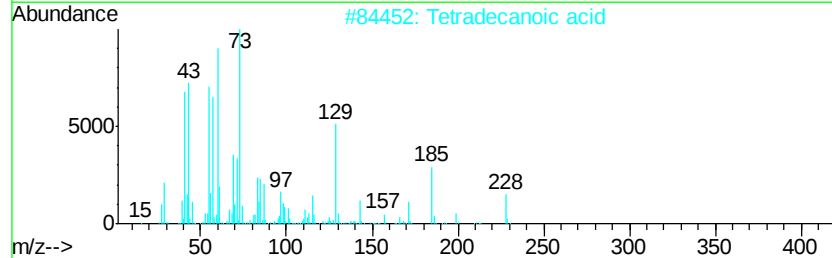
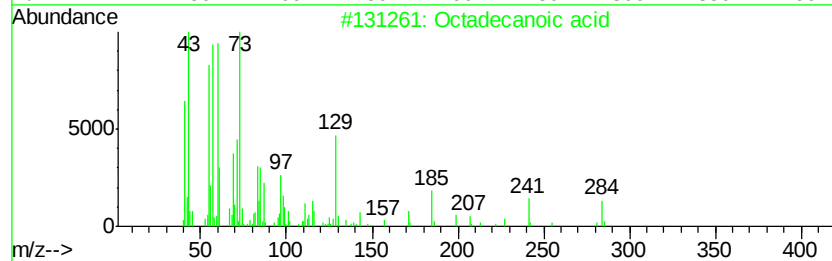
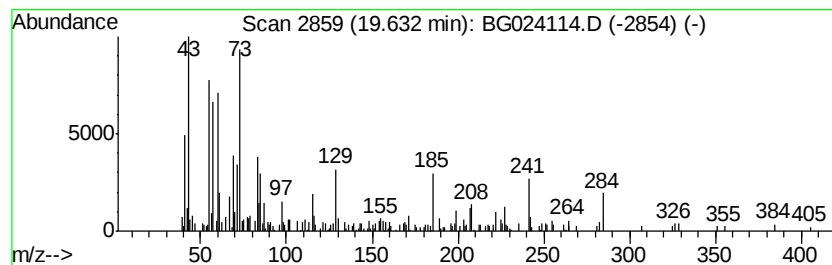
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 6 Octadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.63	3.32 ng	166344	Phenanthrene-d10	17.48

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanoic acid	284	C18H36O2	000057-11-4	94
2	Tetradecanoic acid	228	C14H28O2	000544-63-8	90
3	Pentadecanoic acid	242	C15H30O2	001002-84-2	70
4	n-Decanoic acid	172	C10H20O2	000334-48-5	50
5	Octadecanoic acid, 2-(2-hydroxy...	372	C22H44O4	000106-11-6	47



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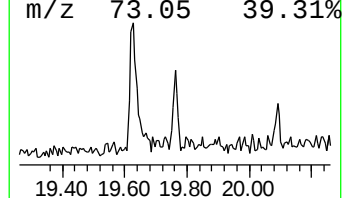
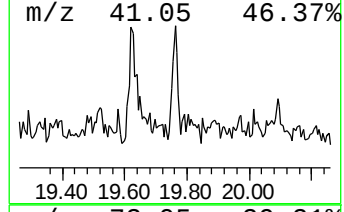
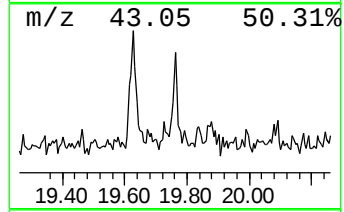
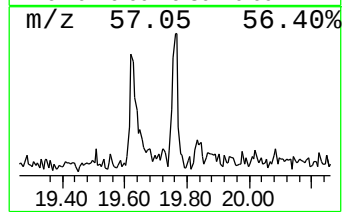
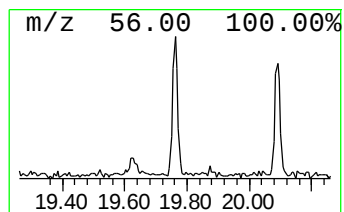
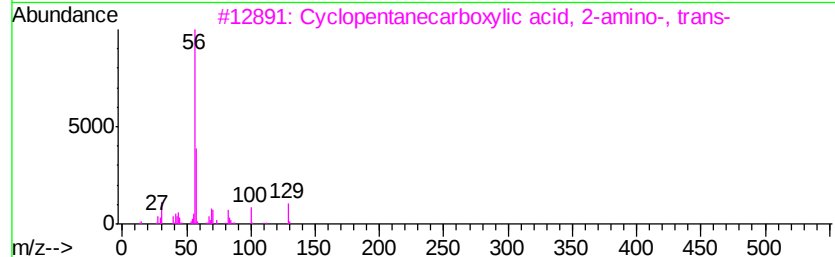
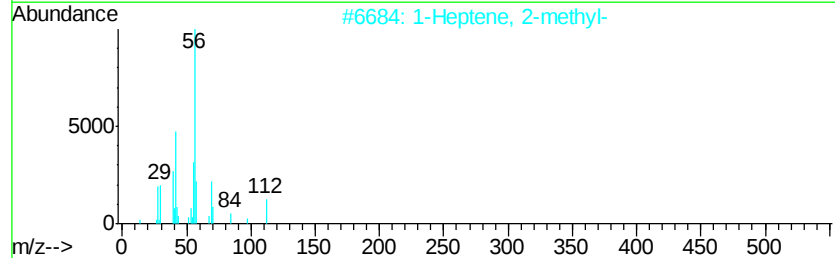
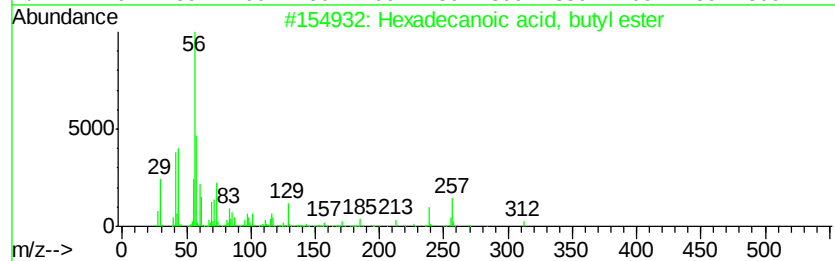
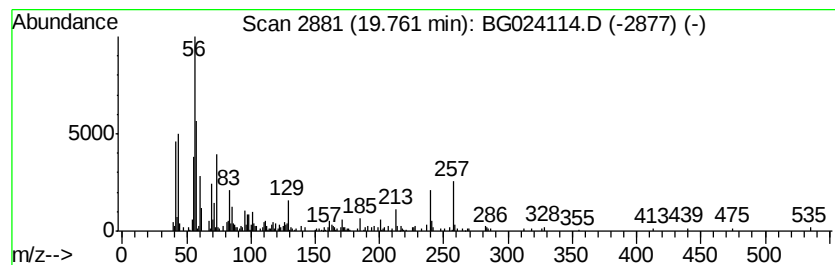
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG092316.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Hexadecanoic acid, butyl ester Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.76	2.12 ng	120821	Chrysene-d12	21.79

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecanoic acid, butyl ester	312	C20H40O2	000111-06-8	59
2		1-Heptene, 2-methyl-	112	C8H16	015870-10-7	43
3		Cyclopentanecarboxylic acid, 2-amino-, trans-	129	C6H11NO2	040482-05-1	38
4		5-Methyl-2-hexanol, chlorodifluoro-	228	C9H15ClF2O2	1000376-27-1	38
5		1-Pentene, 2,4-dimethyl-	98	C7H14	002213-32-3	38



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG092316\
Data File : BG024114.D
Acq On : 24 Sep 2016 8:11
Operator : UM/SJ
Sample : H4893-05
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
PZ-8

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG092316.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.10	6.7	ng	116477	1	8.04	349033	20.0
unknown7.57	7.57	111.9	ng	1952310	1	8.04	349033	20.0
1-Adamantanol	12.12	2.3	ng	63206	2	10.87	550766	20.0
Tridecanoic acid	18.36	2.9	ng	143707	4	17.48	1001770	20.0
Octadecanoic acid	19.63	3.3	ng	166344	4	17.48	1001770	20.0
Hexadecanoic acid...	19.76	2.1	ng	120821	5	21.79	1141240	20.0