

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF082915\
 Data File : BF081277.D
 Acq On : 29 Aug 2015 18:22
 Operator : UM/IZ
 Sample : G3467-06
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW-6

Integration Parameters: rteint.p
 Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0

Filtering: 5
 Min Area: 3 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

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

Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF081715.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	4.560	245	247	257	rVB	304184	431165	16.99%	2.135%
2	5.040	286	289	291	rBV	658155	899734	35.45%	4.455%
3	6.126	382	384	387	rBV	460820	547583	21.57%	2.711%
4	6.240	392	394	397	rBV	1068407	1473384	58.05%	7.295%
5	6.480	413	415	417	rBV	462242	389566	15.35%	1.929%
6	6.640	426	429	431	rBV	1248424	1396879	55.03%	6.917%
7	7.052	462	465	468	rBV	940818	1184329	46.66%	5.864%
8	7.715	519	523	525	rBV	1989373	2538212	100.00%	12.568%
9	7.772	525	528	530	rVB	500291	562401	22.16%	2.785%
10	8.046	550	552	554	rBV	290644	273049	10.76%	1.352%
11	8.366	576	580	582	rBV	22204	26123	1.03%	0.129%
12	8.412	582	584	586	rVB	83342	95216	3.75%	0.471%
13	8.663	603	606	611	rBV2	131675	223418	8.80%	1.106%
14	8.800	611	618	620	rBV	774120	1382837	54.48%	6.847%
15	8.857	620	623	625	rVB	1657750	2120745	83.55%	10.501%
16	9.292	655	661	676	rBV4	128324	514731	20.28%	2.549%
17	9.509	678	680	682	rVB	635083	546000	21.51%	2.704%
18	9.635	688	691	693	rVB	90689	85599	3.37%	0.424%
19	9.795	703	705	707	rBV	35145	47645	1.88%	0.236%
20	9.909	713	715	716	rBV	40618	37828	1.49%	0.187%
21	9.955	716	719	724	rVB2	21292	53748	2.12%	0.266%
22	10.298	746	749	751	rBV	1196713	1226782	48.33%	6.074%
23	10.412	757	759	760	rBV	76047	68437	2.70%	0.339%
24	10.435	760	761	765	rVB	22425	26682	1.05%	0.132%
25	10.983	806	809	810	rBV	417395	547541	21.57%	2.711%
26	11.166	822	825	828	rBV	28364	39522	1.56%	0.196%
27	11.521	854	856	857	rBV	73374	124933	4.92%	0.619%
28	11.543	857	858	860	rVB	142470	108499	4.27%	0.537%
29	11.978	894	896	900	rBV	98214	146750	5.78%	0.727%
30	12.321	924	926	932	rBV	79436	86579	3.41%	0.429%
31	12.572	945	948	950	rBV	1355234	1931969	76.12%	9.566%
32	13.487	1026	1028	1032	rBV	41357	55550	2.19%	0.275%
33	13.589	1035	1037	1040	rBV	562686	519574	20.47%	2.573%
34	14.915	1150	1153	1155	rBV	374699	409717	16.14%	2.029%

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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 >
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35	14.972	1155	1158	1161	rVB	34577	43798	1.73%	0.217%
36	15.692	1218	1221	1225	rBV	17551	29313	1.15%	0.145%

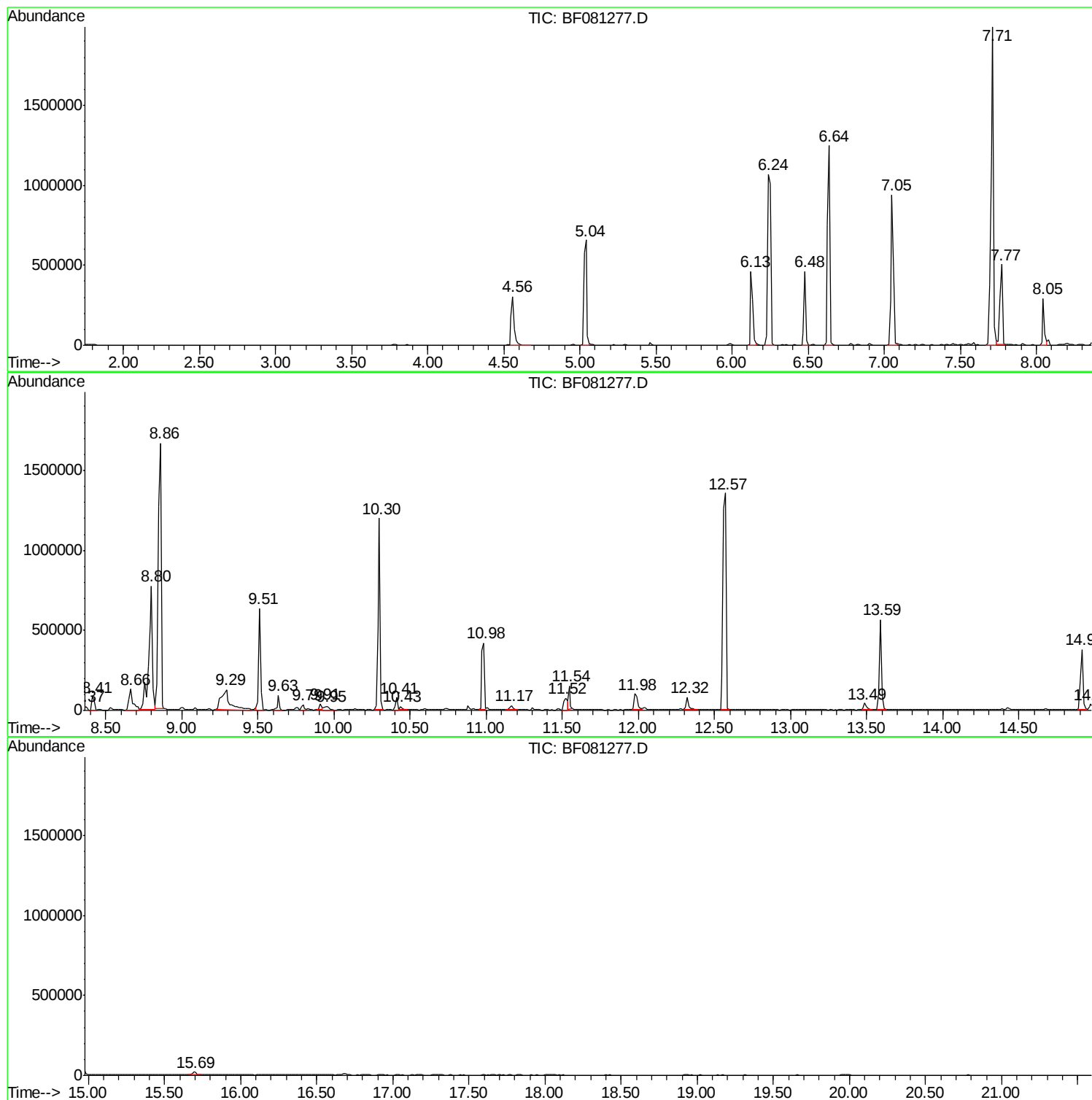
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF081715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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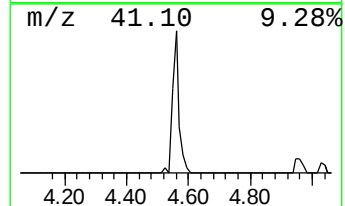
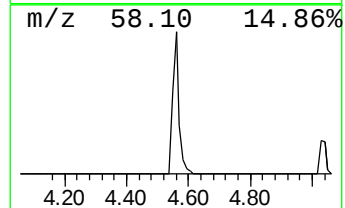
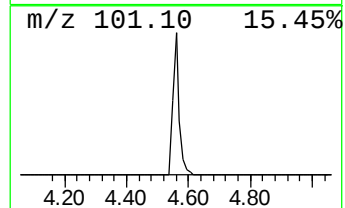
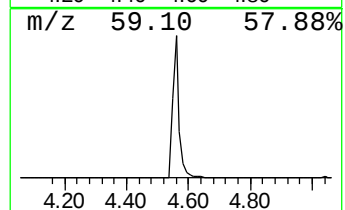
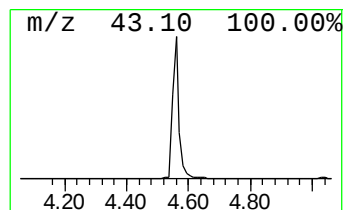
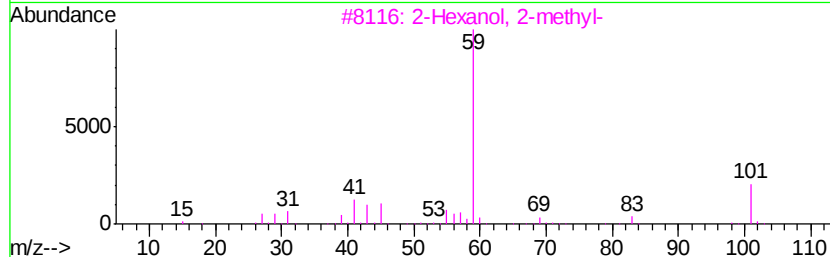
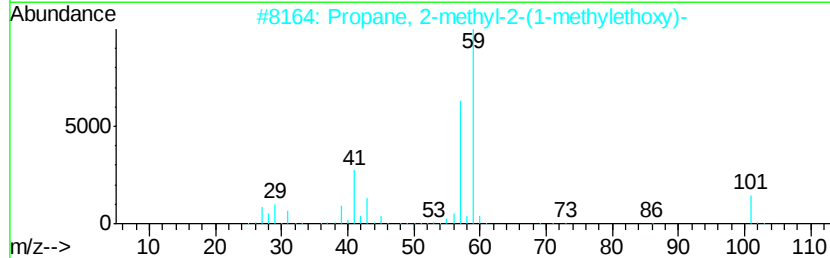
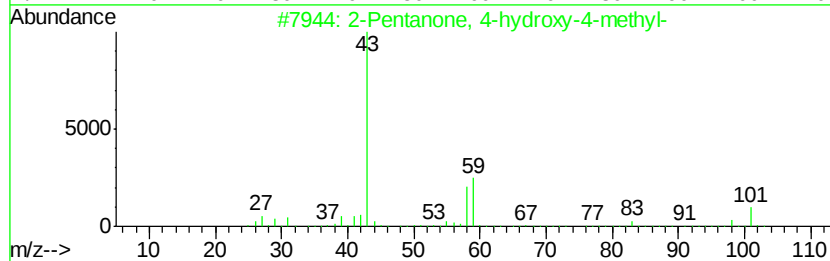
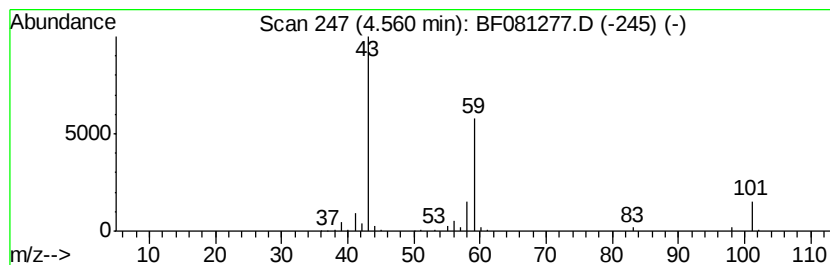
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.56	22.14 ng	431165	1,4-Dichlorobenzene-d4	6.48

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-			116	C6H12O2	000123-42-2	59
2	Propane, 2-methyl-2-(1-methyleth...			116	C7H16O	017348-59-3	36
3	2-Hexanol, 2-methyl-			116	C7H16O	000625-23-0	33
4	3-Hexanol, 4-methyl-			116	C7H16O	000615-29-2	33
5	Acetic acid, cyano-, 1,1-dimethy...			141	C7H11NO2	001116-98-9	32



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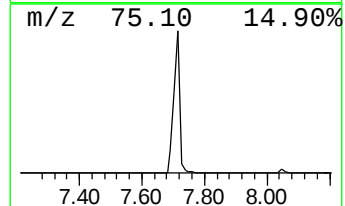
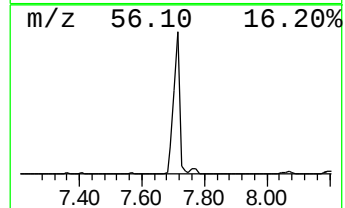
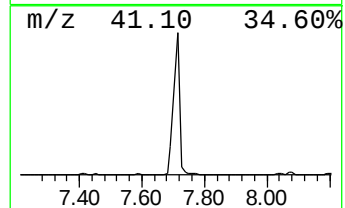
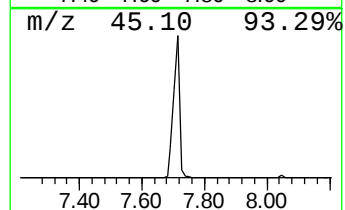
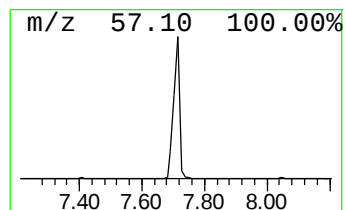
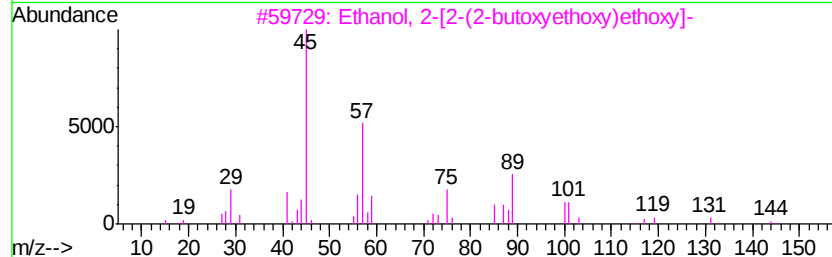
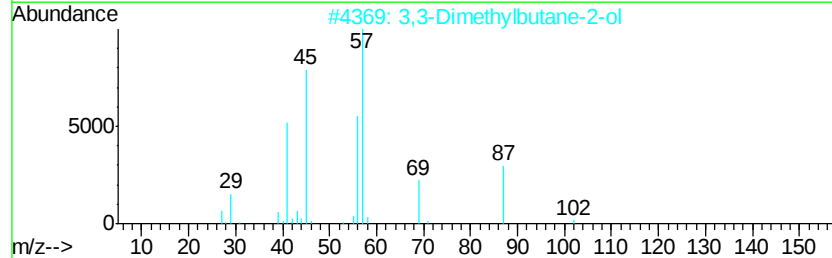
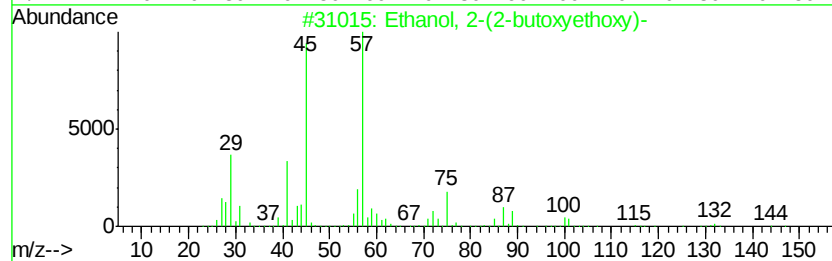
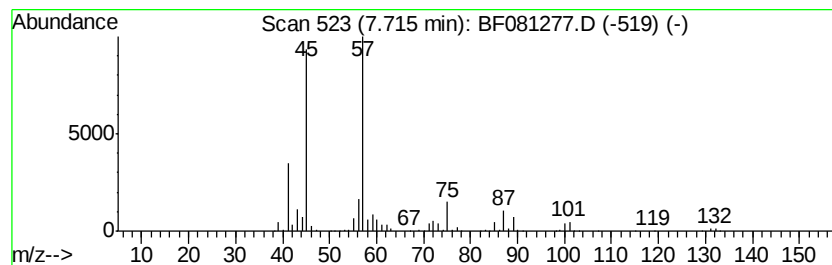
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 4 Ethanol, 2-(2-butoxyethoxy)- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.71	90.26 ng	2538210	Naphthalene-d8	7.77

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
2		3,3-Dimethylbutane-2-ol	102	C6H14O	000464-07-3	53
3		Ethanol, 2-[2-(2-butoxyethoxy)et...	206	C10H22O4	000143-22-6	50
4		Propanal, 3-methoxy-	88	C4H8O2	002806-84-0	47
5		1-Methoxy-3-methyl-3-butene	100	C6H12O	034752-58-4	47



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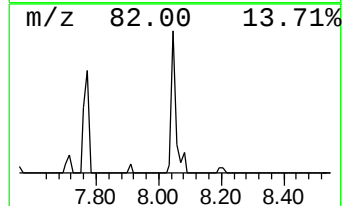
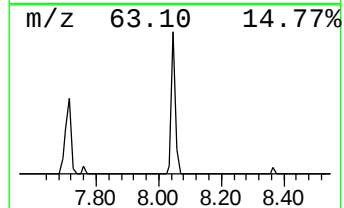
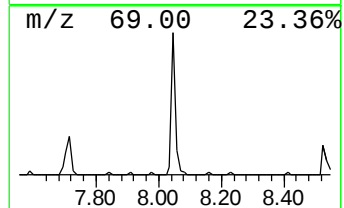
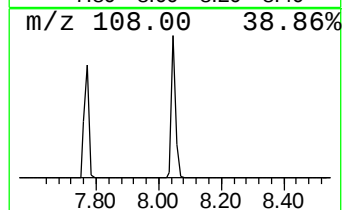
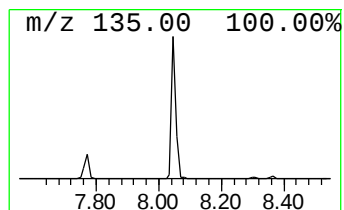
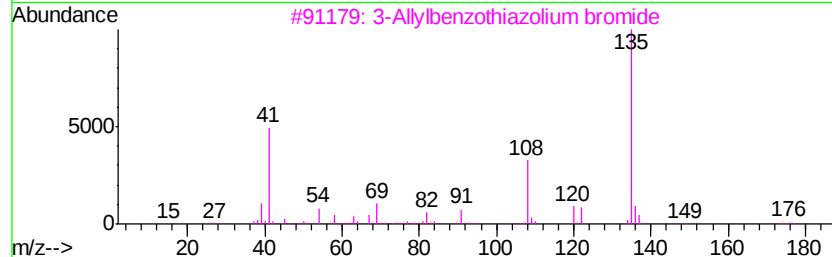
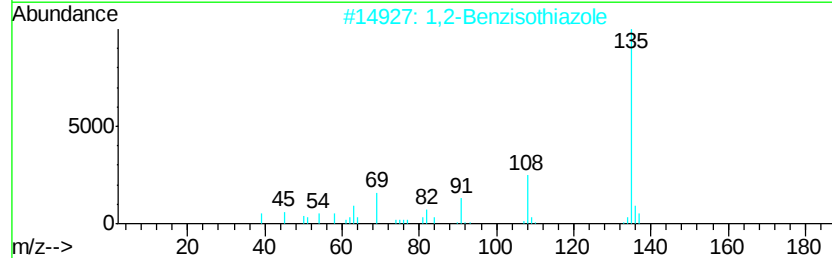
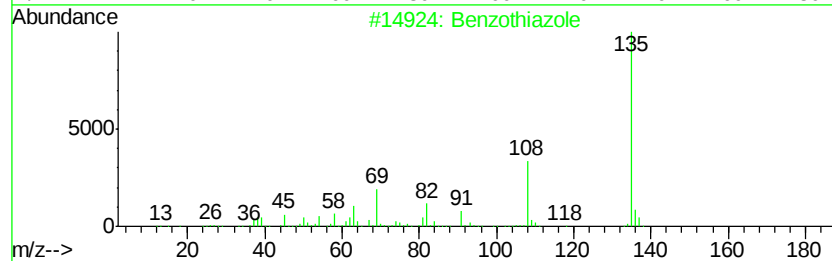
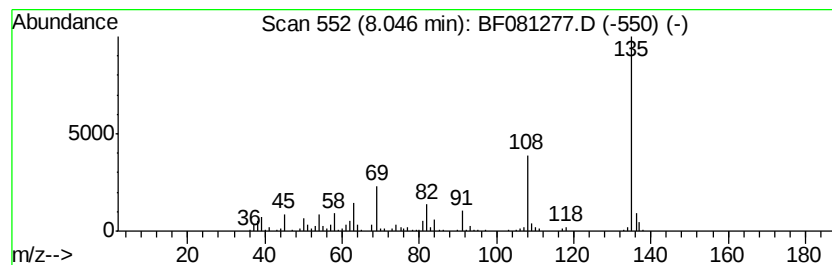
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 5 Benzothiazole Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.05	9.71 ng	273049	Naphthalene-d8	7.77

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzothiazole	135	C7H5NS	000095-16-9	97
2		1,2-Benzisothiazole	135	C7H5NS	000272-16-2	91
3		3-Allylbenzothiazolium bromide	255	C10H10BrNS	016407-55-9	78
4		1H-Pyrazolo[3,4-d]pyrimidin-4-amine	135	C5H5N5	002380-63-4	59
5		4-Aminopyrazolo(3,4-d)pyrimidine	135	C5H5N5	020289-44-5	59



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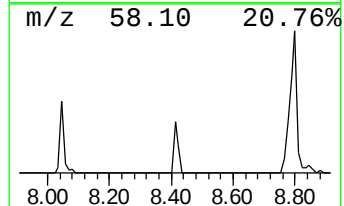
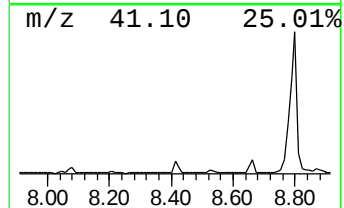
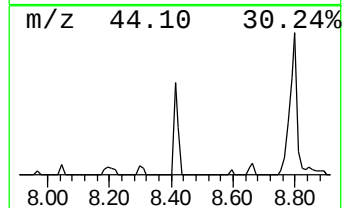
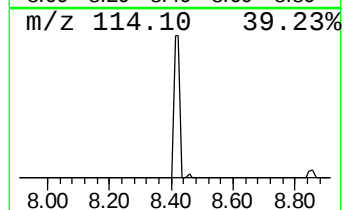
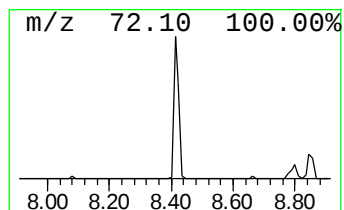
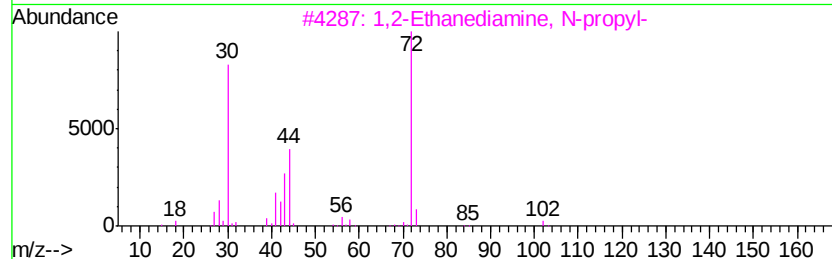
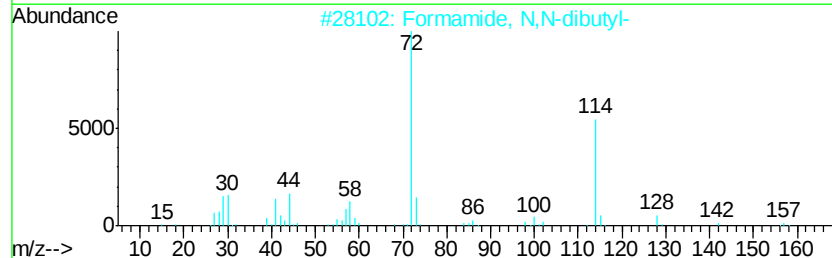
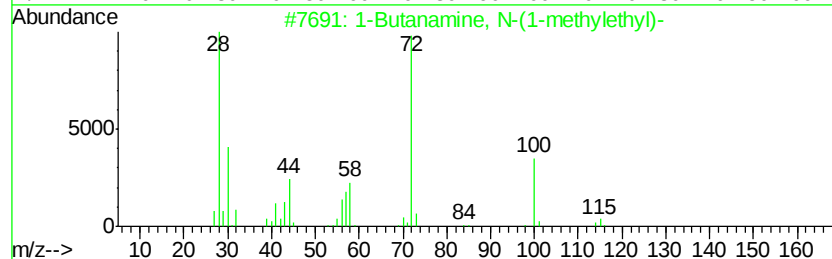
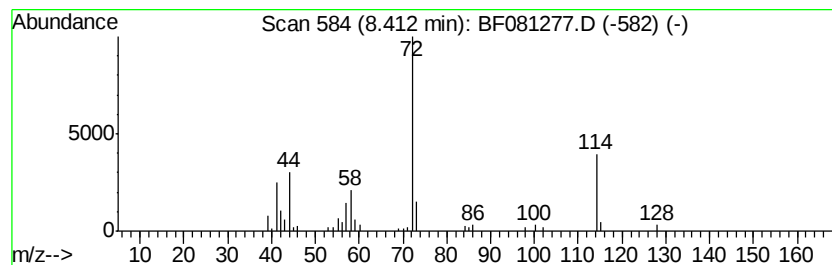
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 1-Butanamine, N-(1-methylet... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.41	3.39 ng	95216	Naphthalene-d8	7.77

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Butanamine, N-(1-methylethyl)-	115	C7H17N	039099-23-5	53	
2	Formamide, N,N-dibutyl-	157	C9H19NO	000761-65-9	50	
3	1,2-Ethanediamine, N-propyl-	102	C5H14N2	000111-39-7	43	
4	1-Butanamine, N-propyl-	115	C7H17N	020193-21-9	38	
5	Isobutyl-propyl-amine	115	C7H17N	039190-66-4	38	



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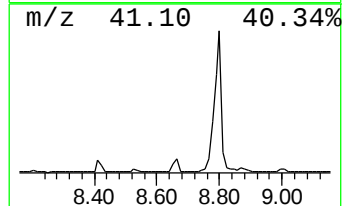
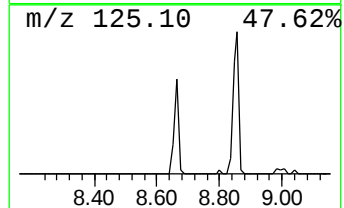
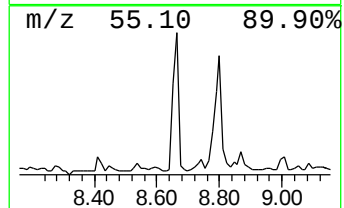
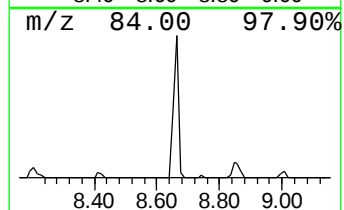
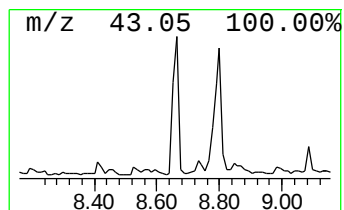
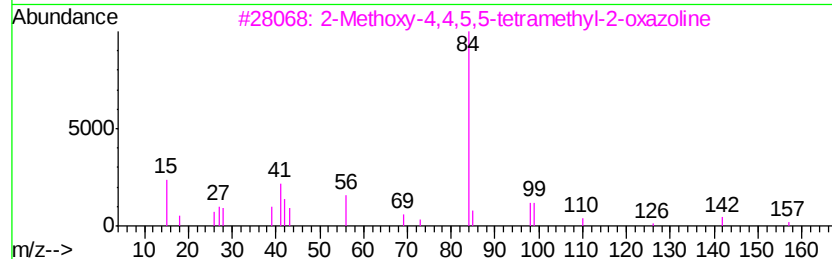
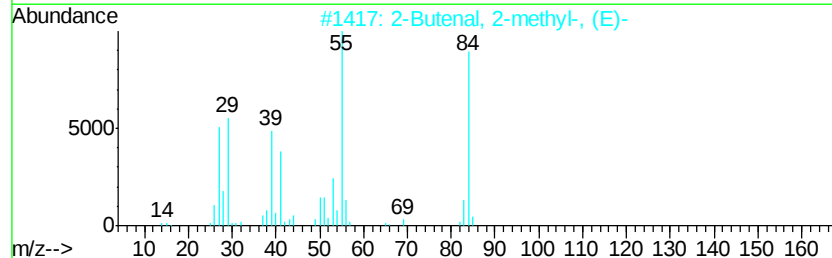
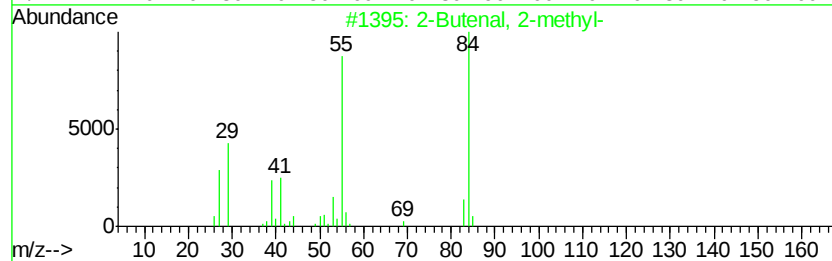
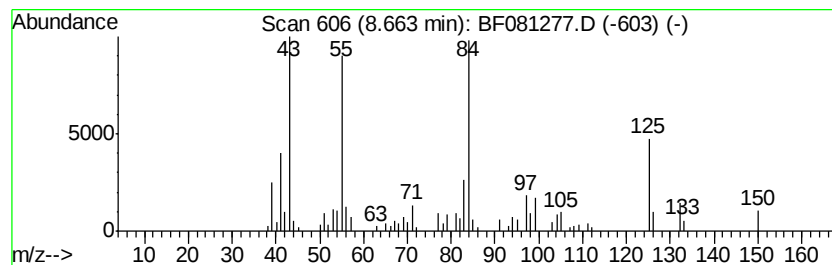
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Peak Number 7 unknown8.66 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.66	8.18 ng	223418	Acenaphthene-d10	9.51

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Butenal, 2-methyl-	84	C5H8O	001115-11-3	43
2		2-Butenal, 2-methyl-, (E)-	84	C5H8O	000497-03-0	38
3		2-Methoxy-4,4,5,5-tetramethyl-2-...	157	C8H15NO2	049680-47-9	35
4		3-Piperidinone, 1-ethyl-	127	C7H13NO	043152-93-8	35
5		1-Pentene, 2,3,3-trimethyl-5-phe...	188	C14H20	1000150-35-4	35



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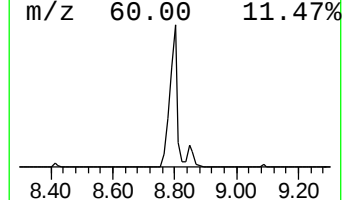
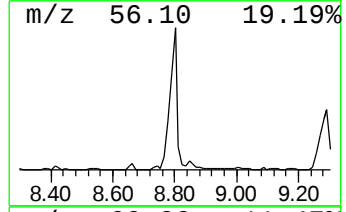
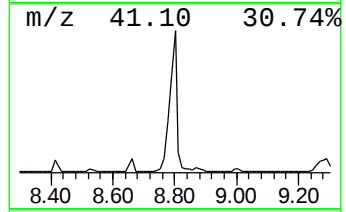
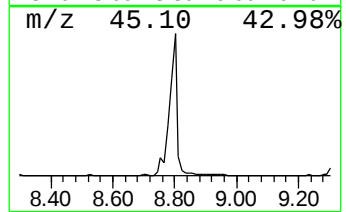
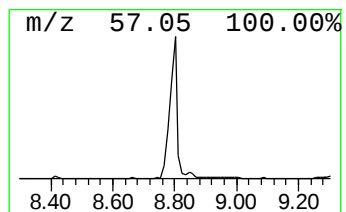
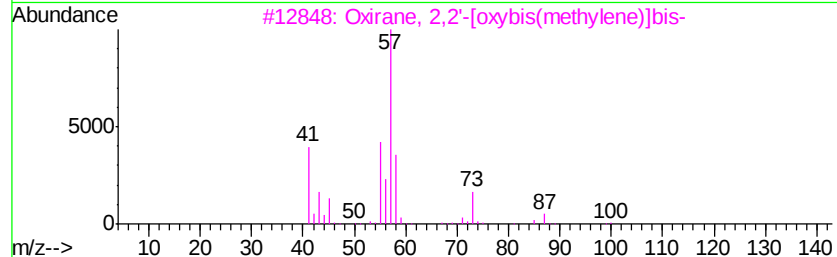
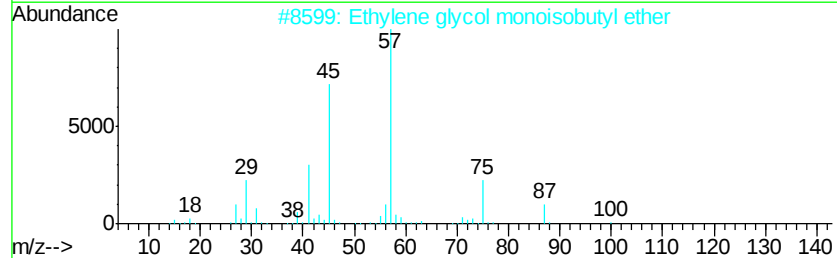
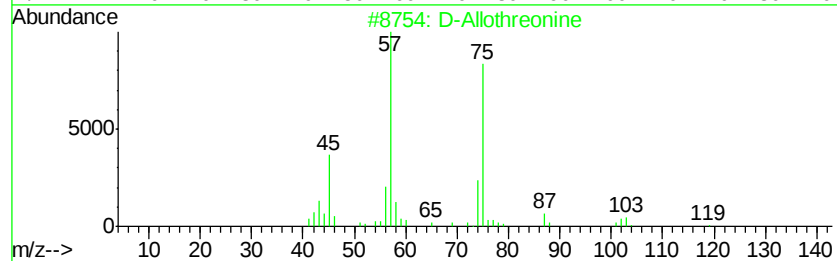
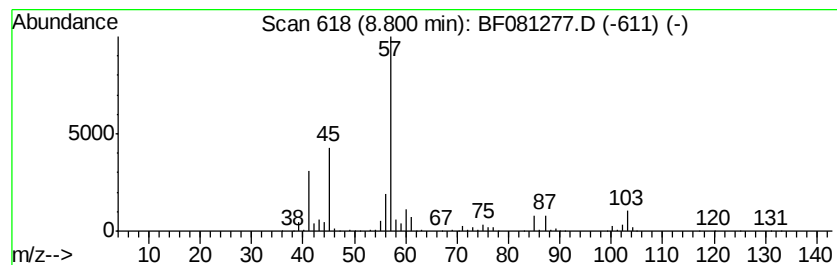
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 unknown8.80 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.80	50.65 ng	1382840	Acenaphthene-d10	9.51

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		D-Allothreonine	119	C4H9NO3	024830-94-2	39
2		Ethylene glycol monoisobutyl ether	118	C6H14O2	004439-24-1	38
3		Oxirane, 2,2'-[oxybis(methylene)]...	130	C6H10O3	002238-07-5	38
4		Ethanol, 2-butoxy-	118	C6H14O2	000111-76-2	37
5		Carbonic acid, bis(1-methylpropy...	174	C9H18O3	000623-63-2	37



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF082915\
Data File : BF081277.D
Acq On : 29 Aug 2015 18:22
Operator : UM/IZ
Sample : G3467-06
Misc :
ALS Vial : 10 Sample Multiplier: 1

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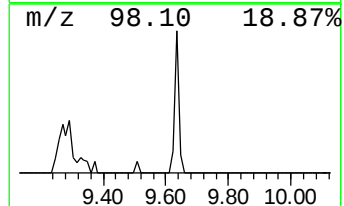
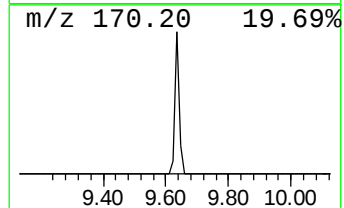
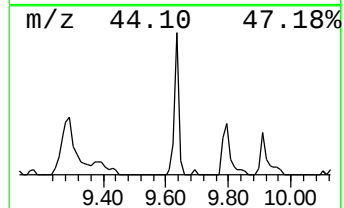
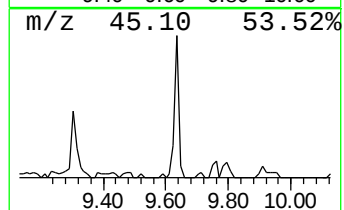
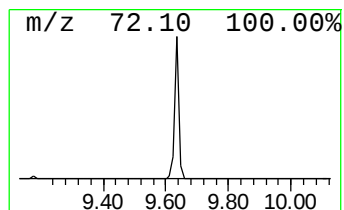
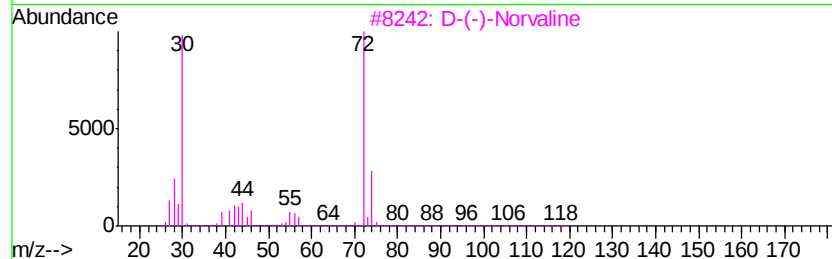
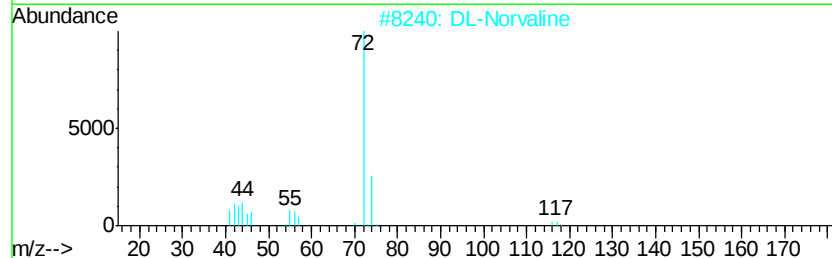
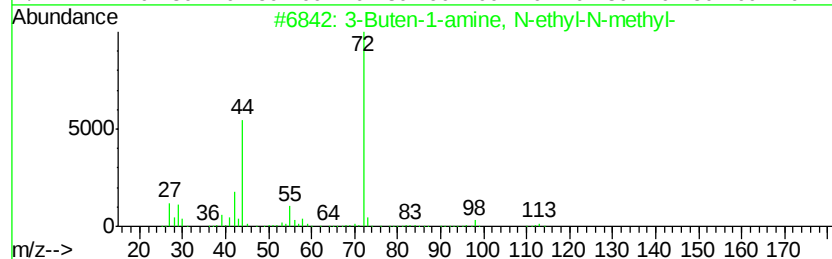
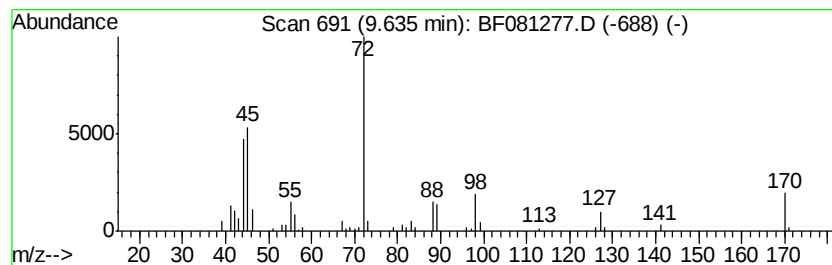
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 unknown9.63 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.63	3.14 ng	85599	Acenaphthene-d10	9.51

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Buten-1-amine, N-ethyl-N-methyl-	113	C7H15N	061308-10-9	38
2			DL-Norvaline	117	C5H11NO2	000760-78-1	25
3			D-(-)-Norvaline	117	C5H11NO2	002013-12-9	25
4			2,3,N,N-Tetramethyl-4-nitro-buty...	188	C8H16N2O3	1000187-66-9	25
5			Ethanamine, N-ethyl-N-methyl-	87	C5H13N	000616-39-7	23



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF082915\
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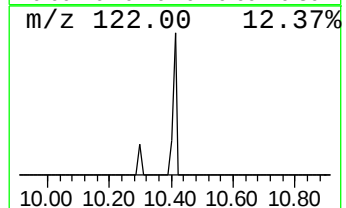
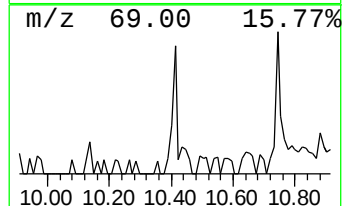
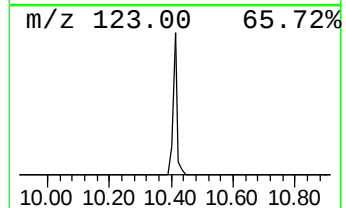
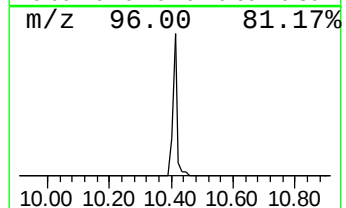
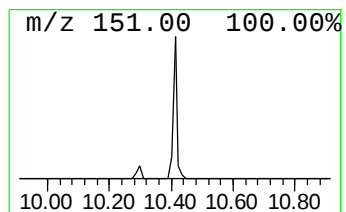
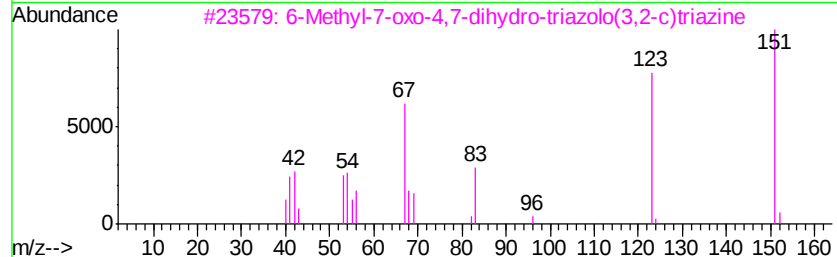
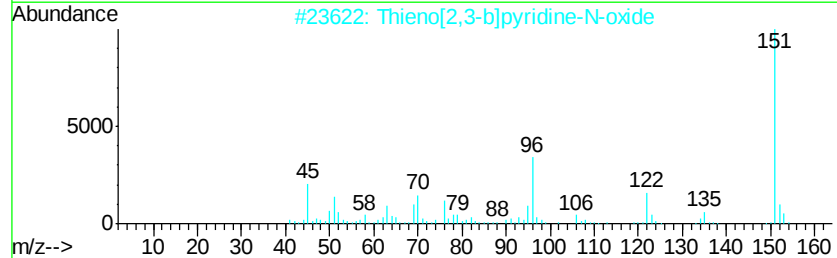
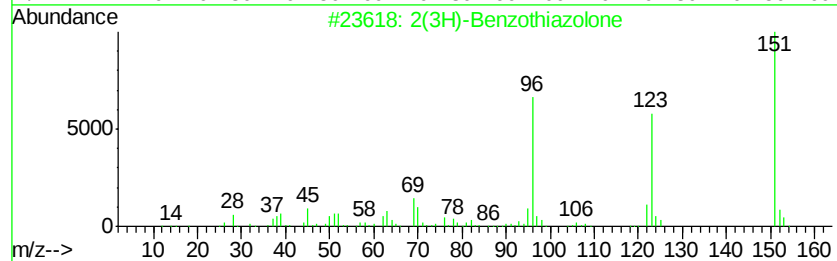
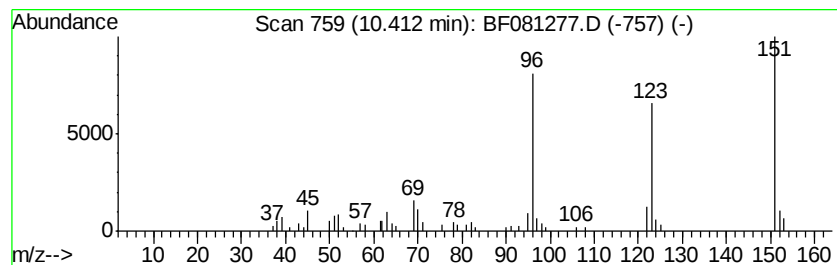
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 2(3H)-Benzothiazolone Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.41	2.50 ng	68437	Phenanthrene-d10	10.98	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2(3H)-Benzothiazolone	151	C7H5NOS	000934-34-9	95
2	Thieno[2,3-b]pyridine-N-oxide	151	C7H5NOS	025557-50-0	46
3	6-Methyl-7-oxo-4,7-dihydro-triaz...	151	C5H5N5O	057250-39-2	38
4	4-Methoxyformanilide	151	C8H9NO2	023896-88-0	35
5	1,2-Benzisothiazol-3(2H)-one	151	C7H5NOS	002634-33-5	35



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF082915\
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Acq On : 29 Aug 2015 18:22
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Sample : G3467-06
Misc :
ALS Vial : 10 Sample Multiplier: 1

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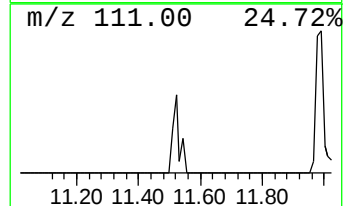
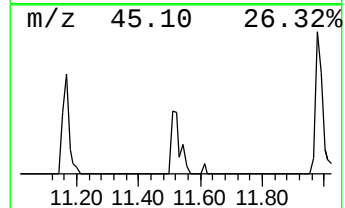
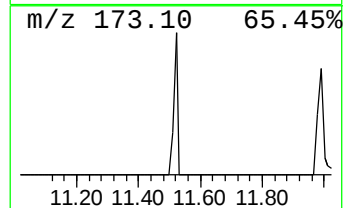
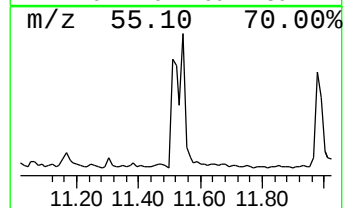
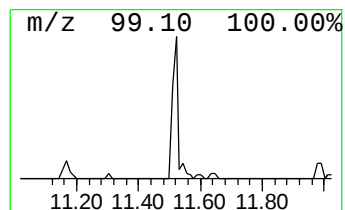
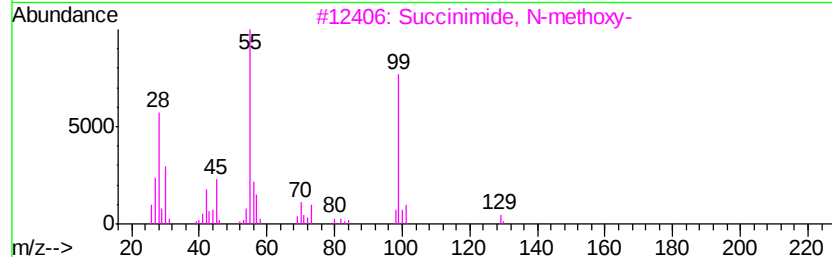
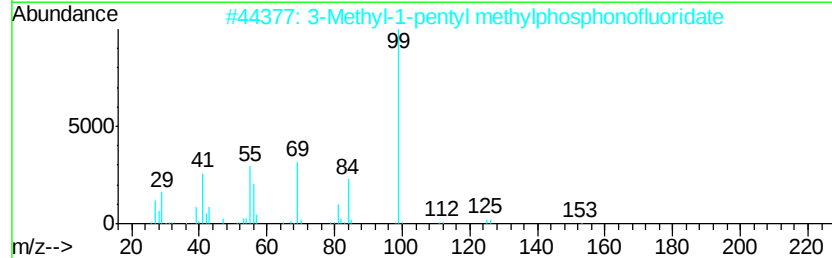
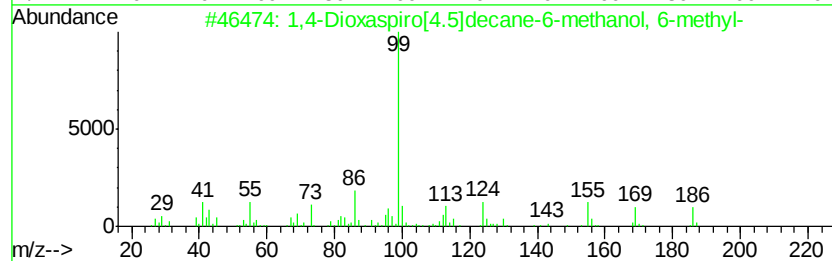
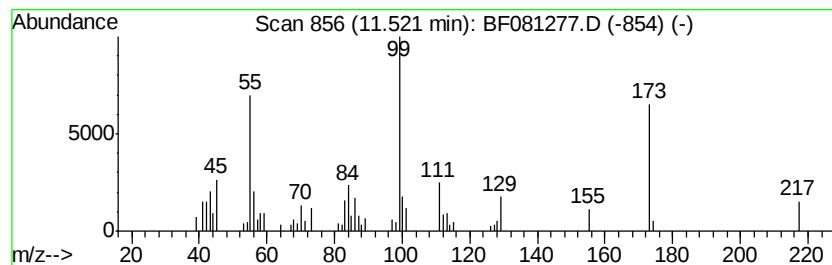
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 unknown11.52 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.52	4.56 ng	124933	Phenanthrene-d10	10.98

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,4-Dioxaspiro[4.5]decane-6-meth...	186	C10H18O3	087787-09-5	35
2			3-Methyl-1-pentyl methylphosphon...	182	C7H16F02P	199850-60-7	35
3			Succinimide, N-methoxy-	129	C5H7N03	005904-50-7	25
4			Methylphosphonic acid, fluoroanh...	238	C11H24F02P	211192-71-1	25
5			4H-Cyclohepta[b]thiophen-4-one, ...	294	C16H26N2OS	1000265-44-1	25



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF082915\
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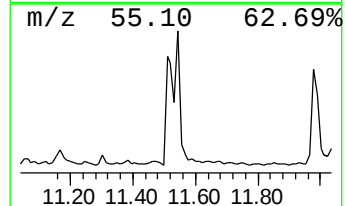
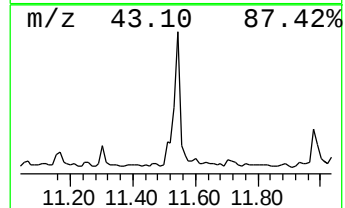
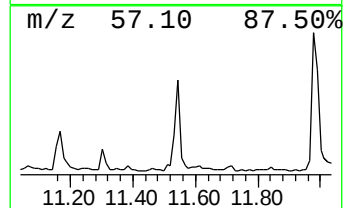
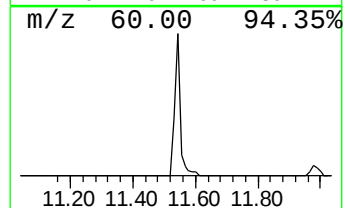
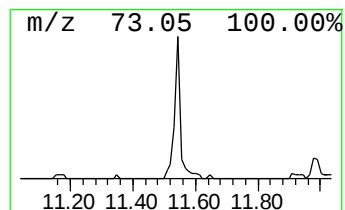
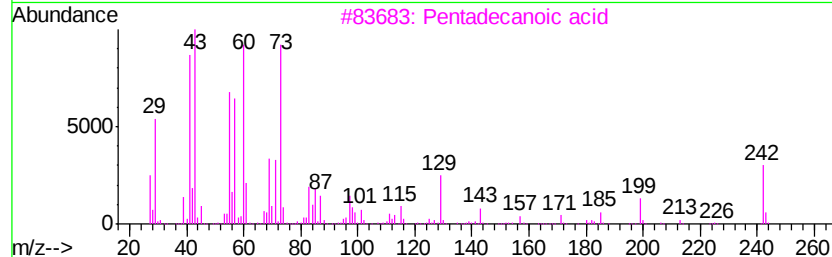
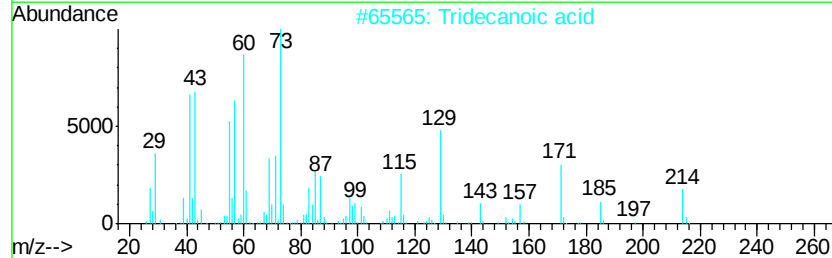
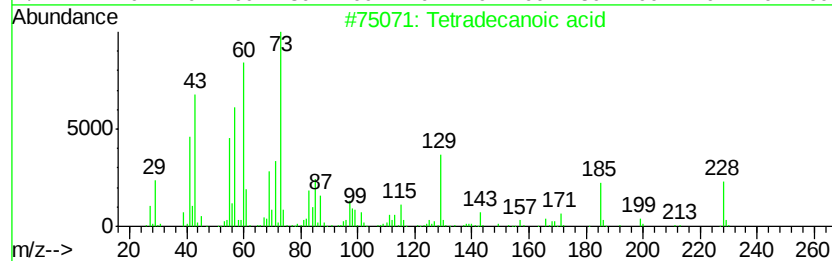
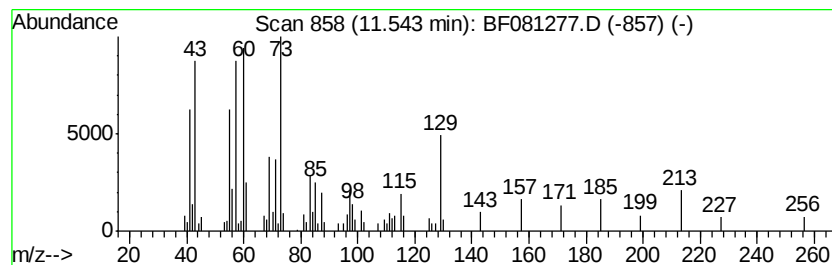
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Tetradecanoic acid Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.54	3.96 ng	108499	Phenanthrene-d10	10.98

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetradecanoic acid	228	C14H28O2	000544-63-8	80
2			Tridecanoic acid	214	C13H26O2	000638-53-9	64
3			Pentadecanoic acid	242	C15H30O2	001002-84-2	64
4			n-Decanoic acid	172	C10H20O2	000334-48-5	52
5			Dodecanoic acid	200	C12H24O2	000143-07-7	50



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF082915\
Data File : BF081277.D
Acq On : 29 Aug 2015 18:22
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Misc :
ALS Vial : 10 Sample Multiplier: 1

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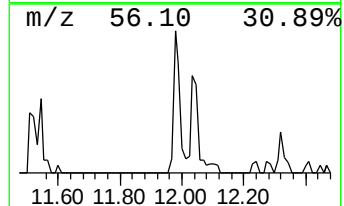
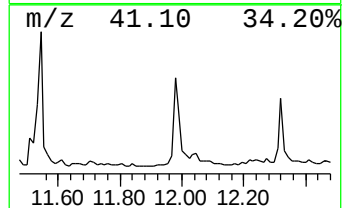
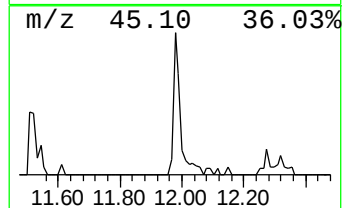
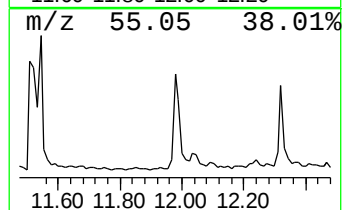
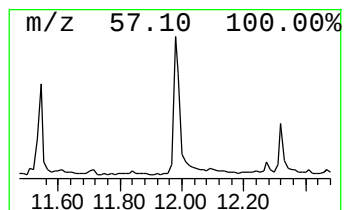
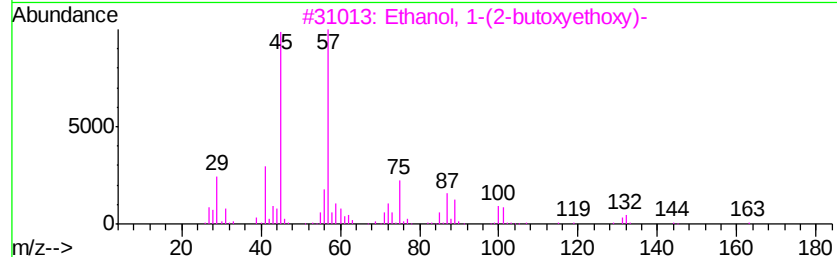
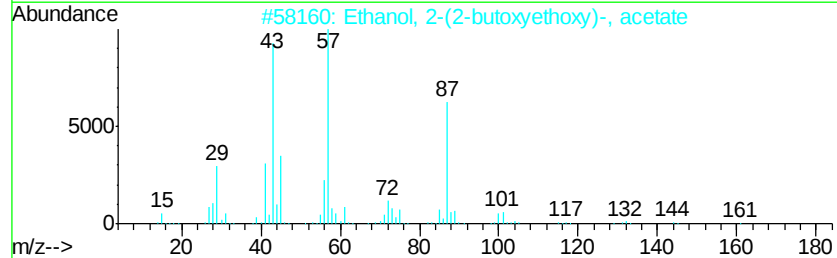
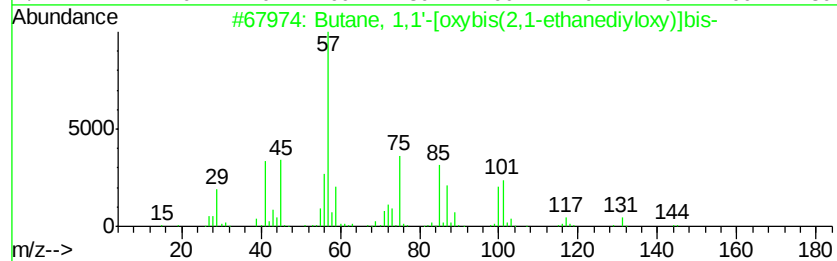
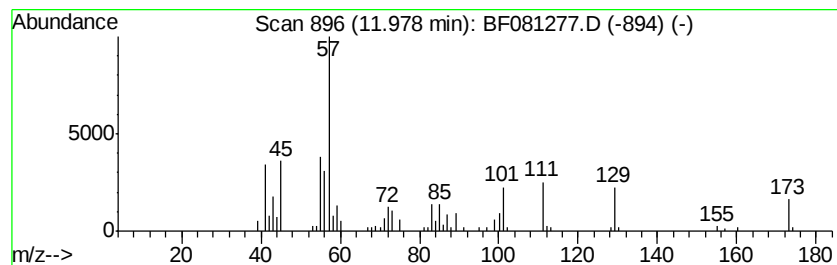
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 unknown11.98 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.98	5.36 ng	146750	Phenanthrene-d10	10.98

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 1,1'-[oxybis(2,1-ethaned...	218	C12H26O3	000112-73-2	27
2		Ethanol, 2-(2-butoxyethoxy)-, ac...	204	C10H20O4	000124-17-4	25
3		Ethanol, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	12
4		6-Dodecanone, 5,8-diethyl-7-hydr...	256	C16H32O2	031814-59-2	12
5		Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	10



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF082915\
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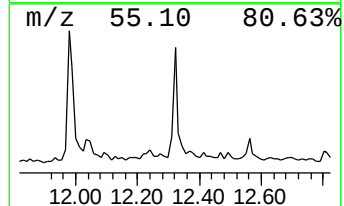
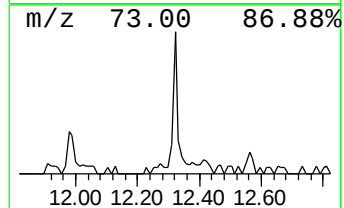
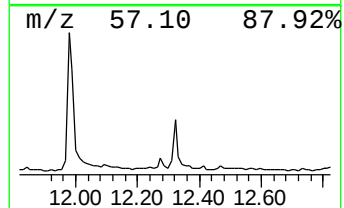
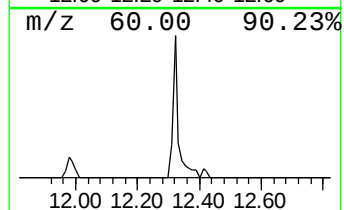
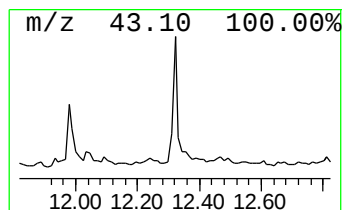
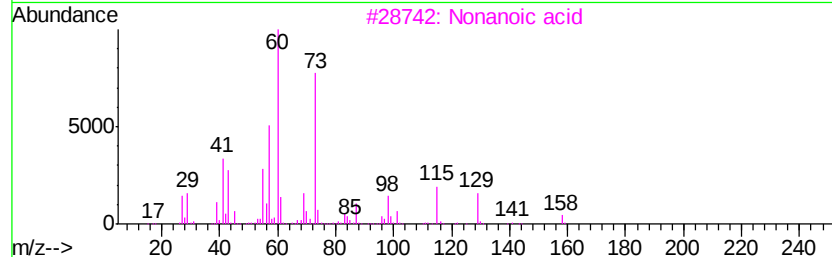
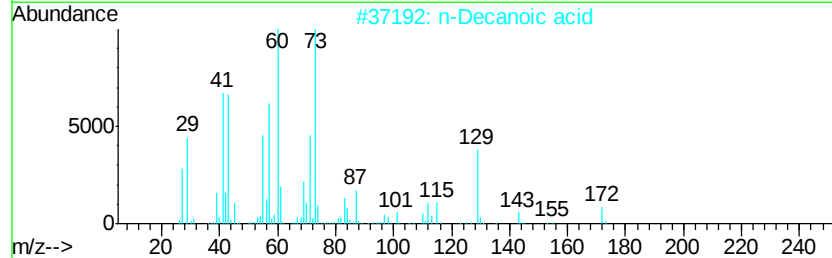
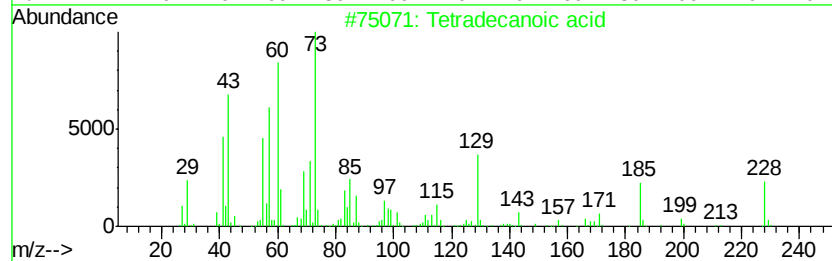
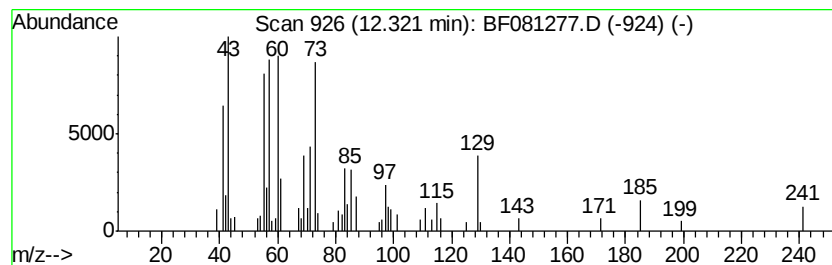
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 n-Decanoic acid Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.32	3.33 ng	86579	Chrysene-d12	13.59
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Tetradecanoic acid	228 C14H28O2	000544-63-8 72
2		n-Decanoic acid	172 C10H20O2	000334-48-5 53
3		Nonanoic acid	158 C9H18O2	000112-05-0 35
4		Amyl Nitrite	117 C5H11NO2	000110-46-3 35
5		1-Octanol, 2-butyl-	186 C12H26O	003913-02-8 14



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF082915\
Data File : BF081277.D
Acq On : 29 Aug 2015 18:22
Operator : UM/IZ
Sample : G3467-06
Misc :
ALS Vial : 10 Sample Multiplier: 1

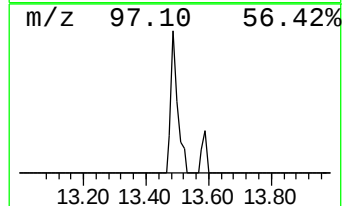
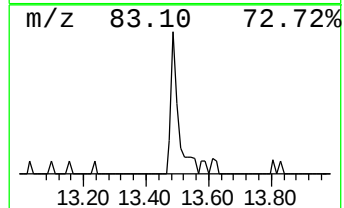
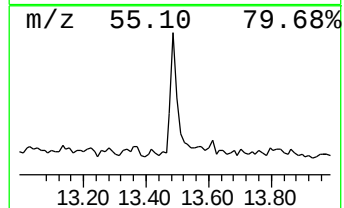
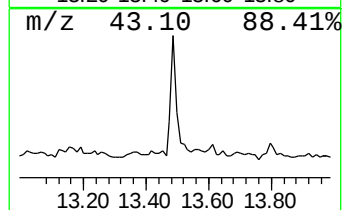
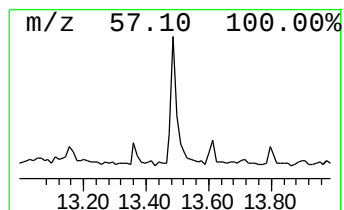
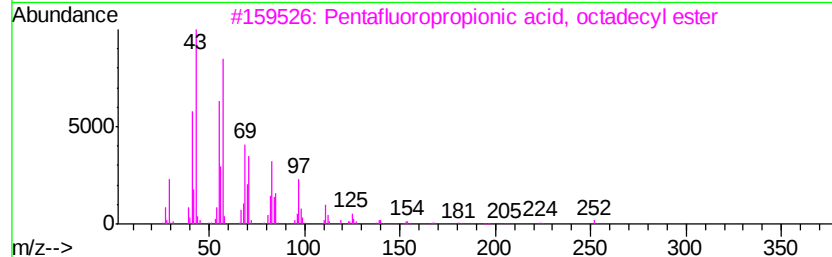
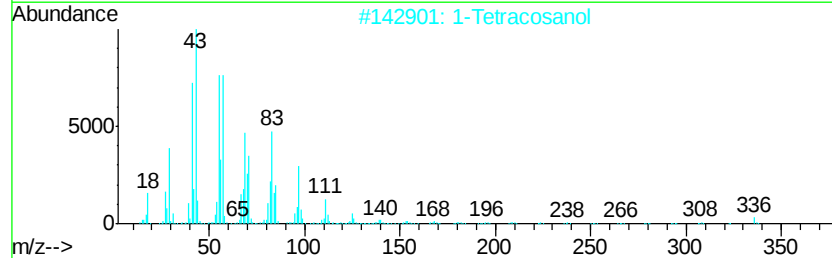
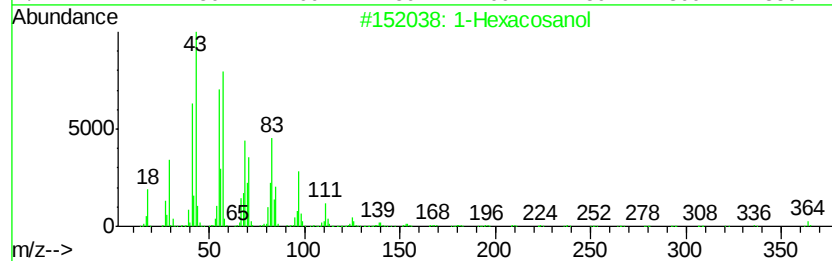
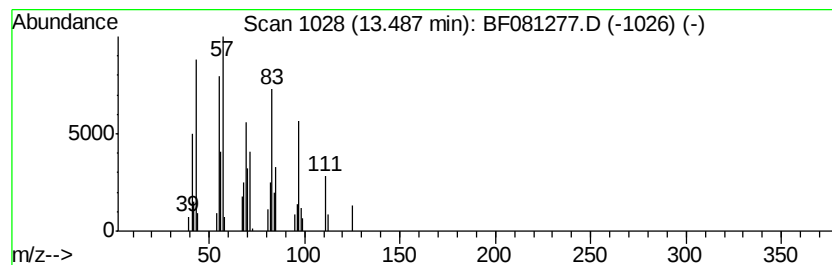
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF081715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 1-Hexacosanol Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.49	2.14 ng	55550	Chrysene-d12	13.59
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	1-Hexacosanol	382 C26H54O	000506-52-5	87
2	1-Tetracosanol	354 C24H50O	000506-51-4	74
3	Pentafluoropropionic acid, octad...	416 C21H37F5O2	1000280-07-7	72
4	Octadecane, 1-(ethenyloxy)-	296 C20H40O	000930-02-9	64
5	Silane, trichlorodocosyl-	442 C22H45Cl3Si	007325-84-0	59



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF082915\
Data File : BF081277.D
Acq On : 29 Aug 2015 18:22
Operator : UM/IZ
Sample : G3467-06
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MW-6

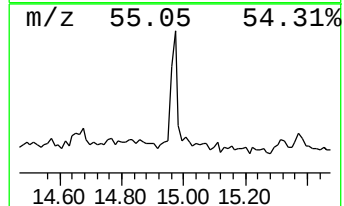
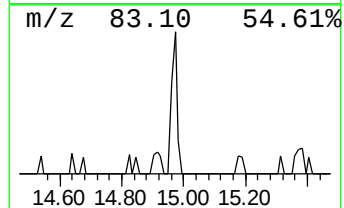
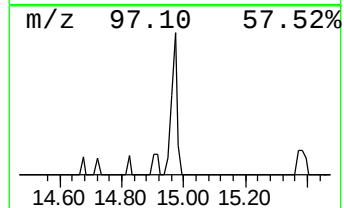
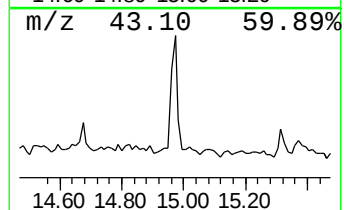
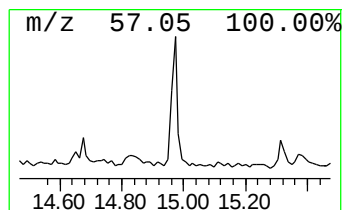
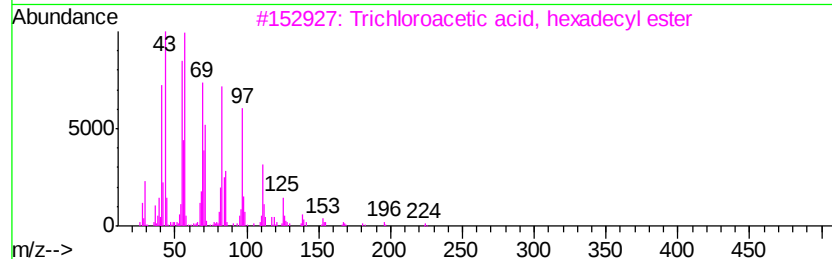
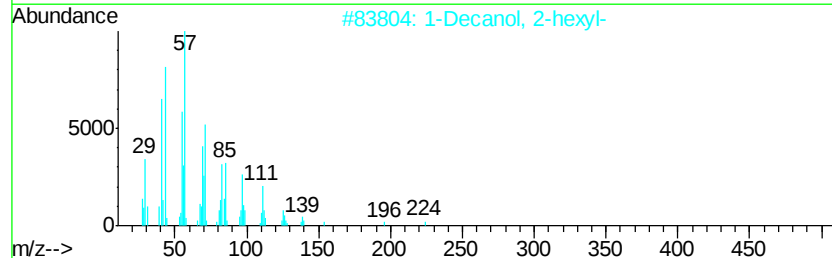
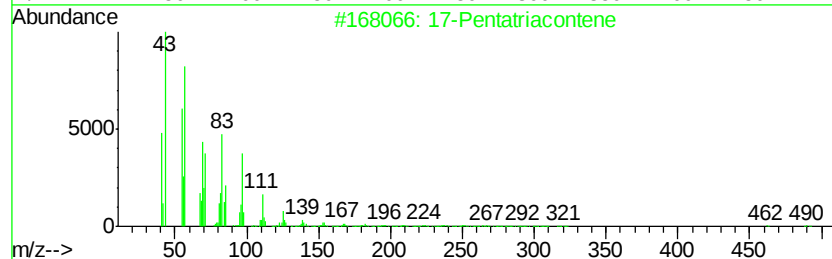
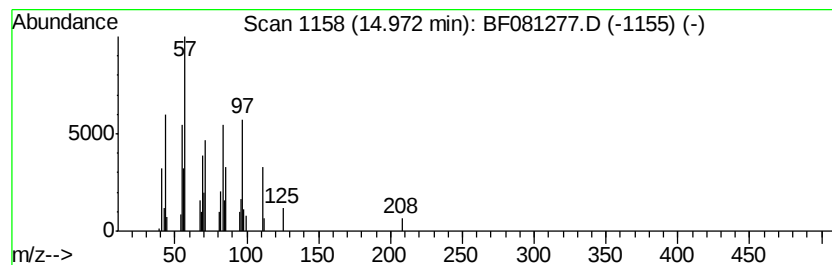
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF081715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 17-Pentatriacontene Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.97	2.14 ng	43798	Perylene-d12	14.92

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			17-Pentatriacontene	491	C35H70	006971-40-0	58
2			1-Decanol, 2-hexyl-	242	C16H34O	002425-77-6	50
3			Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	46
4			Trichloroacetic acid, pentadecyl...	372	C17H31Cl3O2	074339-53-0	46
5			Trichloroacetic acid, tetradecyl...	358	C16H29Cl3O2	074339-52-9	43



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 Data File : BF081277.D
 Acq On : 29 Aug 2015 18:22
 Operator : UM/IZ
 Sample : G3467-06
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW-6

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF081715.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
2-Pentanone, 4-hy...	4.56	22.1	ng	431165	1	6.48	389566	20.0
Ethanol, 2-(2-but...	7.71	90.3	ng	2538210	2	7.77	562401	20.0
Benzothiazole	8.05	9.7	ng	273049	2	7.77	562401	20.0
1-Butanamine, N-(...	8.41	3.4	ng	95216	2	7.77	562401	20.0
unknown8.66	8.66	8.2	ng	223418	3	9.51	546000	20.0
unknown8.80	8.80	50.6	ng	1382840	3	9.51	546000	20.0
unknown9.63	9.63	3.1	ng	85599	3	9.51	546000	20.0
2(3H)-Benzothiazo...	10.41	2.5	ng	68437	4	10.98	547541	20.0
unknown11.52	11.52	4.6	ng	124933	4	10.98	547541	20.0
Tetradecanoic acid	11.54	4.0	ng	108499	4	10.98	547541	20.0
unknown11.98	11.98	5.4	ng	146750	4	10.98	547541	20.0
n-Decanoic acid	12.32	3.3	ng	86579	5	13.59	519574	20.0
1-Hexacosanol	13.49	2.1	ng	55550	5	13.59	519574	20.0
17-Pentatriacontene	14.97	2.1	ng	43798	6	14.92	409717	20.0