

Data Path : Z:\HPCHEM1\BNA F\DATA\BF033115\
 Data File : BF078121.D
 Acq On : 31 Mar 2015 18:37
 Operator : TP/IZ
 Sample : G1700-01
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 S8-0424

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-BF031115.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	1.321	35	38	41	rVB	241590	353261	18.92%	2.575%
2	1.401	43	45	49	rVV2	12493	21154	1.13%	0.154%
3	1.481	49	52	55	rVB	325505	385975	20.68%	2.813%
4	1.618	62	64	71	rVV	49941	108700	5.82%	0.792%
5	1.721	71	73	111	rVB	771649	1369791	73.38%	9.985%
6	2.338	125	127	132	rVB	10019	19271	1.03%	0.140%
7	4.773	339	340	347	rBB	26347	40157	2.15%	0.293%
8	5.333	386	389	394	rBV	384731	479335	25.68%	3.494%
9	6.636	501	503	509	rBV	288307	297343	15.93%	2.167%
10	6.762	511	514	516	rBV	1026262	995079	53.31%	7.253%
11	7.047	537	539	541	rBV	225821	237311	12.71%	1.730%
12	7.230	552	555	557	rBV	1084714	973827	52.17%	7.098%
13	7.333	562	564	569	rVV	13589	23091	1.24%	0.168%
14	7.745	597	600	602	rBV	811240	798722	42.79%	5.822%
15	7.802	603	605	607	rVB	19800	19809	1.06%	0.144%
16	8.110	627	632	635	rVB	41422	49830	2.67%	0.363%
17	8.625	675	677	679	rBV	289702	344960	18.48%	2.514%
18	8.819	691	694	699	rBV4	10905	20174	1.08%	0.147%
19	9.162	722	724	728	rBV3	12218	20686	1.11%	0.151%
20	9.276	732	734	737	rVB	22477	24939	1.34%	0.182%
21	9.550	756	758	762	rBV4	14016	25291	1.35%	0.184%
22	9.676	768	769	775	rVB4	32269	68061	3.65%	0.496%
23	9.813	775	781	786	rBV6	16947	56454	3.02%	0.411%
24	9.905	786	789	792	rBV4	17199	39628	2.12%	0.289%
25	9.973	792	795	797	rVV	1694175	1724217	92.37%	12.568%
26	10.019	797	799	803	rBV4	11001	23376	1.25%	0.170%
27	10.191	812	814	817	rVB2	27186	41071	2.20%	0.299%
28	10.339	824	827	829	rBV2	25885	42395	2.27%	0.309%
29	10.396	829	832	833	rVB2	12947	25663	1.37%	0.187%
30	10.476	837	839	844	rVB5	27348	34029	1.82%	0.248%
31	10.556	844	846	848	rBV	40023	53196	2.85%	0.388%
32	10.625	850	852	855	rVB4	11008	22457	1.20%	0.164%
33	10.693	855	858	862	rBV4	10450	23290	1.25%	0.170%
34	10.785	864	866	869	rBV	404706	422325	22.62%	3.078%

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Integration Parameters: rteint.p
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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

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Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	11.048	886	889	891 rBV	18271	28557	1.53%	0.208%
36	11.105	893	894	896 rVV	12689	19103	1.02%	0.139%
37	11.151	896	898	902 rVV2	29639	78369	4.20%	0.571%
38	11.231	902	905	910 rVB3	65283	107054	5.73%	0.780%
39	11.699	943	946	948 rVB2	14648	24757	1.33%	0.180%
40	11.768	950	952	955 rBV	969744	976367	52.30%	7.117%
41	12.202	988	990	993 rVB2	18262	25884	1.39%	0.189%
42	12.465	1012	1013	1016 rVB2	23223	28565	1.53%	0.208%
43	12.534	1016	1019	1020 rBV3	9768	21463	1.15%	0.156%
44	12.625	1024	1027	1029 rVB	461392	451612	24.19%	3.292%
45	14.625	1199	1202	1204 rBV	1513216	1866691	100.00%	13.607%
46	15.802	1303	1305	1311 rBV3	16106	32318	1.73%	0.236%
47	15.894	1311	1313	1316 rBV	329286	457403	24.50%	3.334%
48	17.631	1462	1465	1469 rBV	309805	416076	22.29%	3.033%

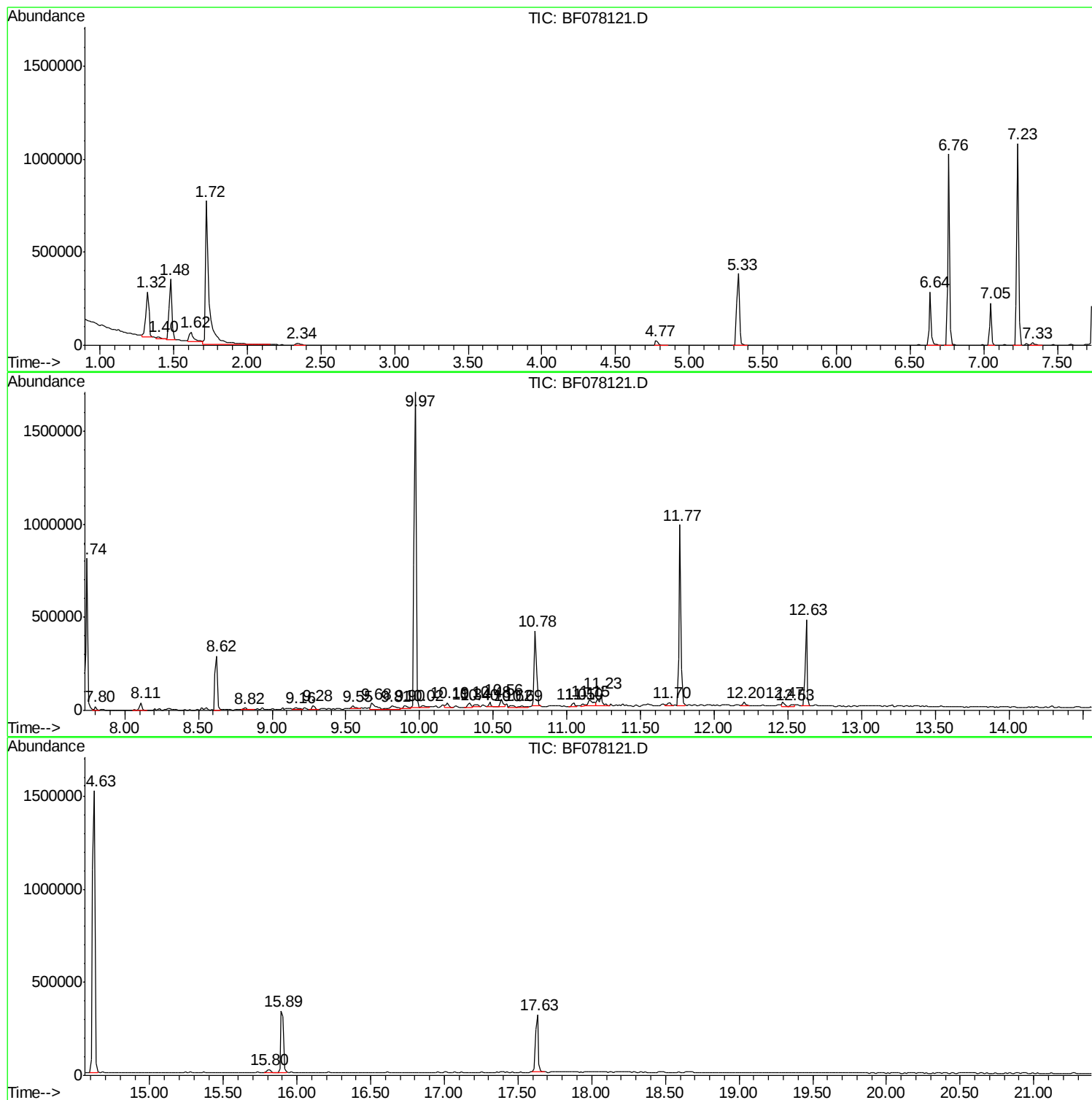
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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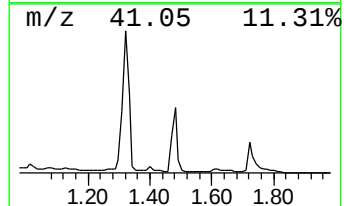
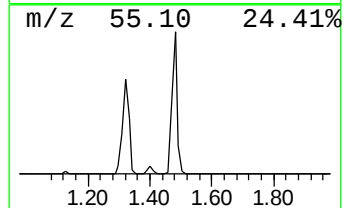
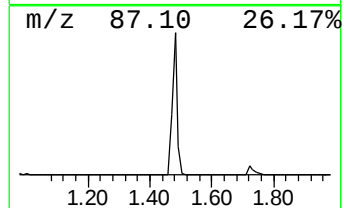
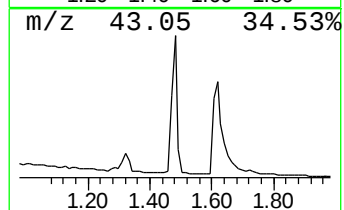
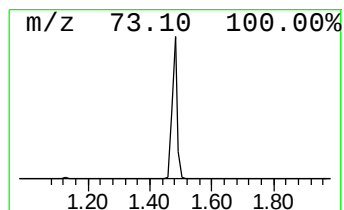
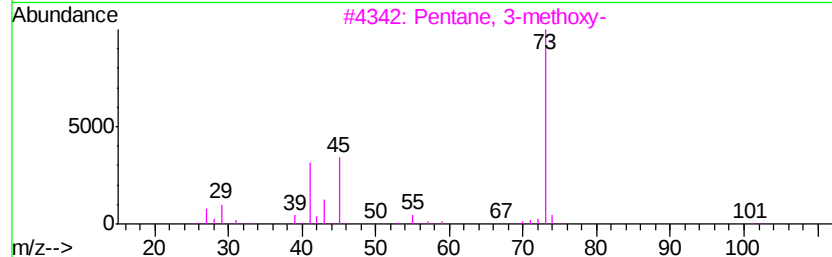
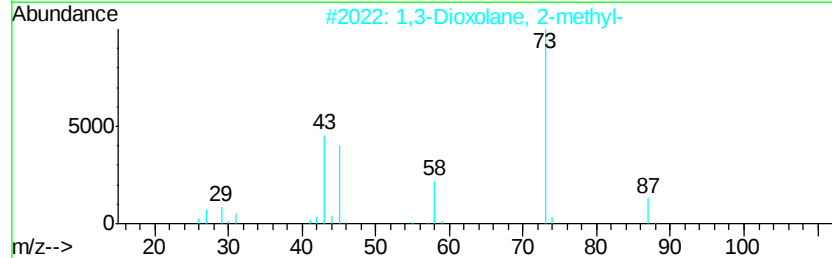
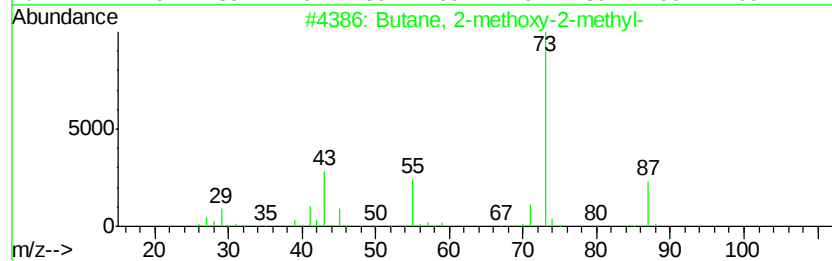
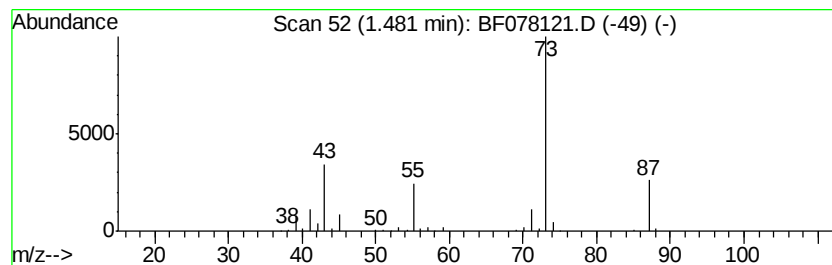
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 Butane, 2-methoxy-2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.48	32.53 ng	385975	1,4-Dichlorobenzene-d4	7.05

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2		1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	37
3		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	23
4		Silane, tetramethyl-	88	C4H12Si	000075-76-3	17
5		Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	9



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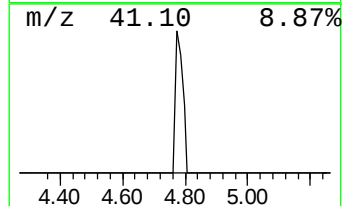
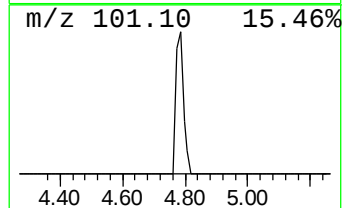
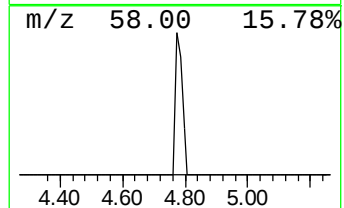
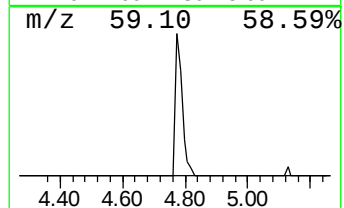
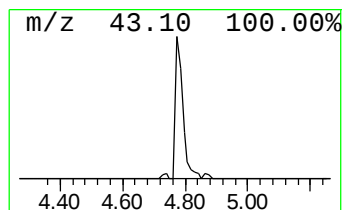
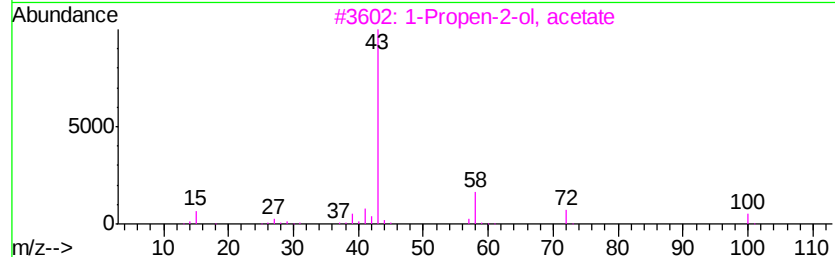
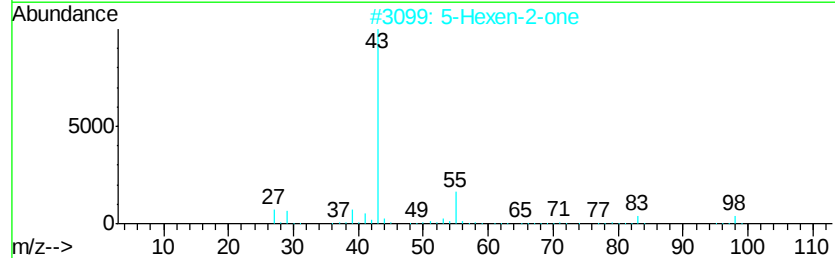
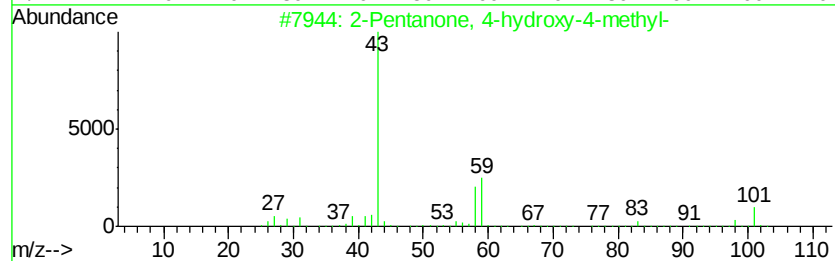
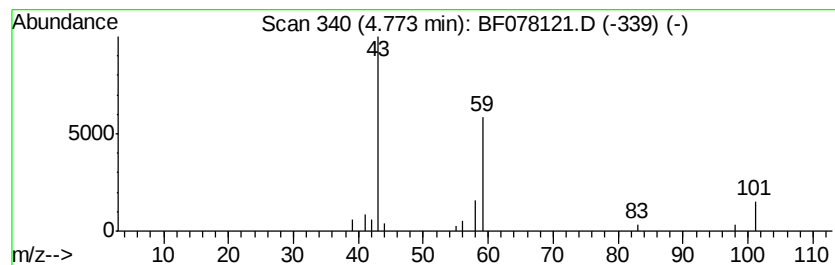
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.77	3.38 ng	40157	1,4-Dichlorobenzene-d4	7.05

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2		5-Hexen-2-one	98	C6H10O	000109-49-9	9
3		1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	9
4		2-Propanone, 1-(1-methylethoxy)-	116	C6H12O2	042781-12-4	9
5		Diacetamide	101	C4H7NO2	000625-77-4	7



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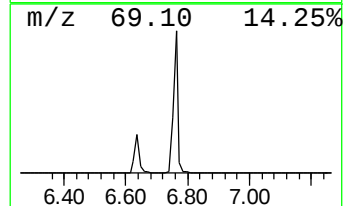
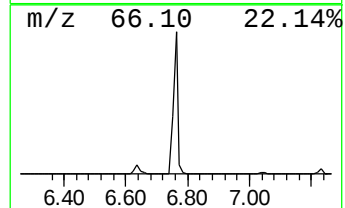
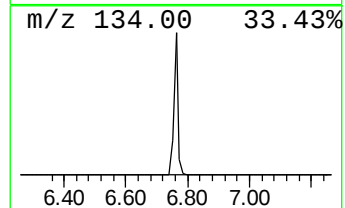
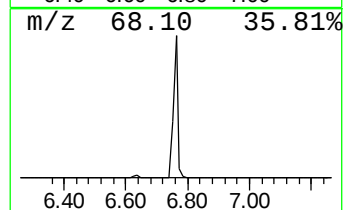
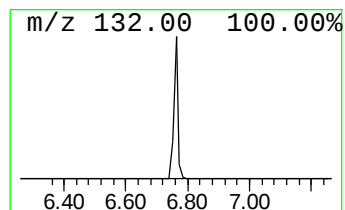
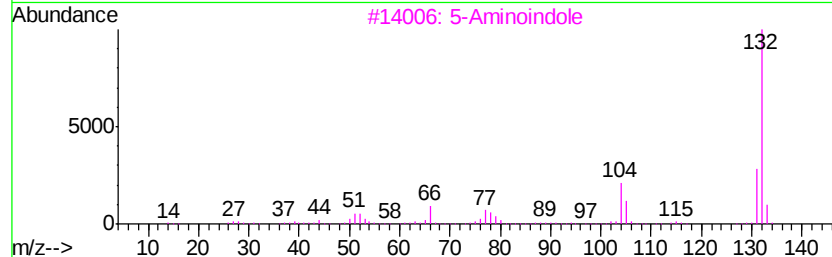
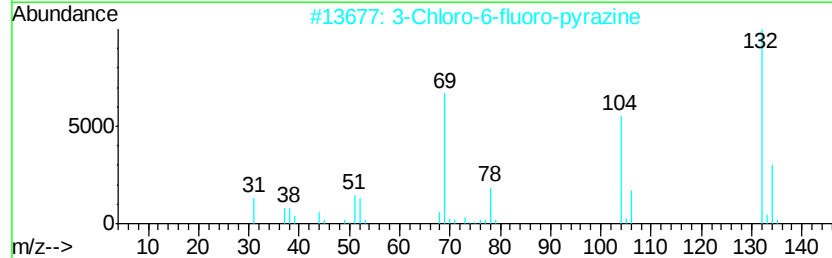
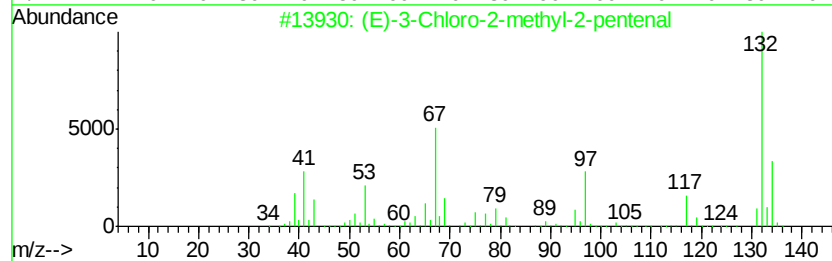
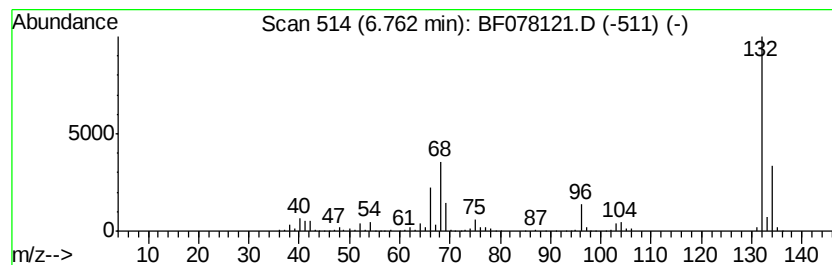
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 unknown6.76 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.76	83.86 ng	995079	1,4-Dichlorobenzene-d4	7.05

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	25
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	25
3		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9
5		4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	9



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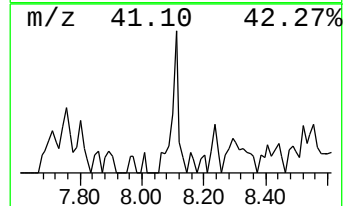
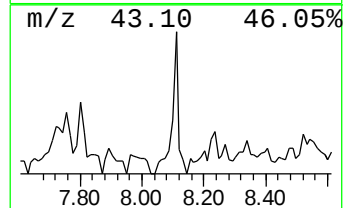
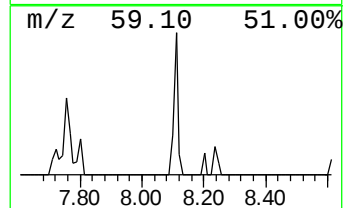
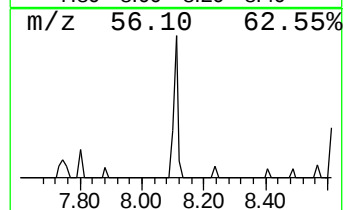
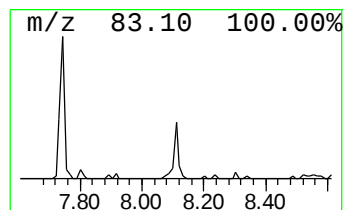
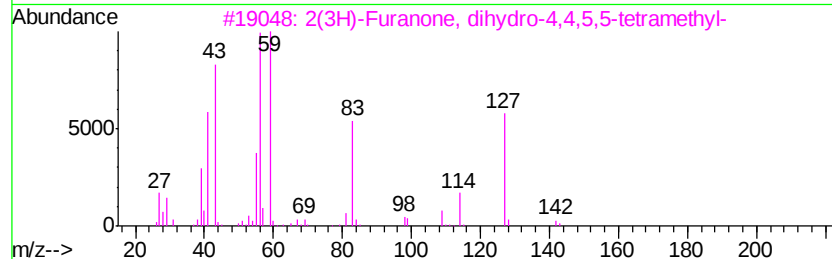
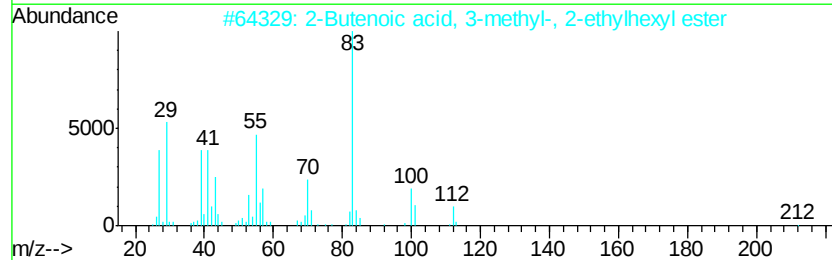
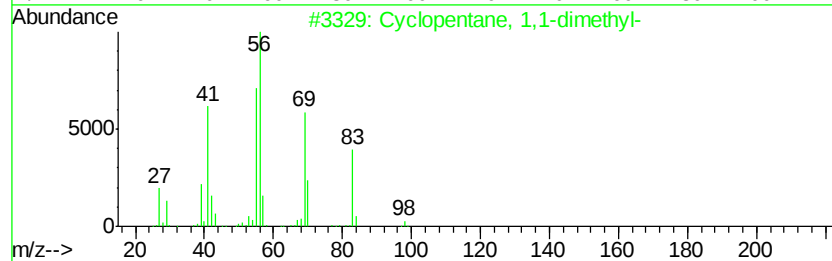
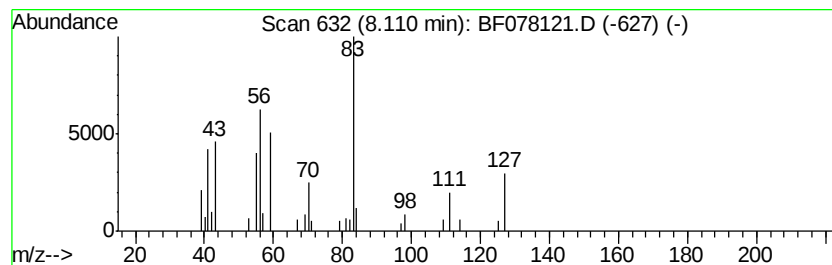
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 unknown8.11 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.11	2.89 ng	49830	Napthalene-d8	8.62

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopentane, 1,1-dimethyl-	98	C7H14	001638-26-2	49
2		2-Butenoic acid, 3-methyl-, 2-ethylhexyl ester	212	C13H24O2	085567-31-3	35
3		2(3H)-Furanone, dihydro-4,4,5,5-tetramethyl-	142	C8H14O2	016466-24-3	28
4		3-Nonen-5-one	140	C9H16O	082456-34-6	17
5		Cyclopentane, 2-ethyl-1,1-dimethyl-	126	C9H18	054549-80-3	12



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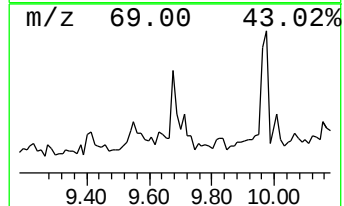
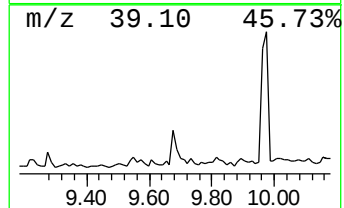
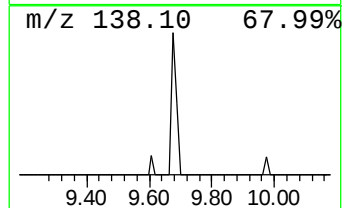
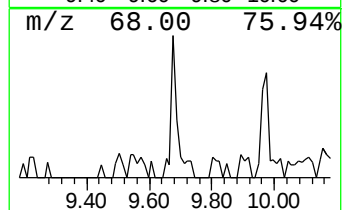
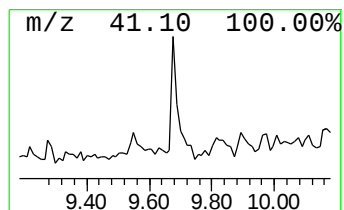
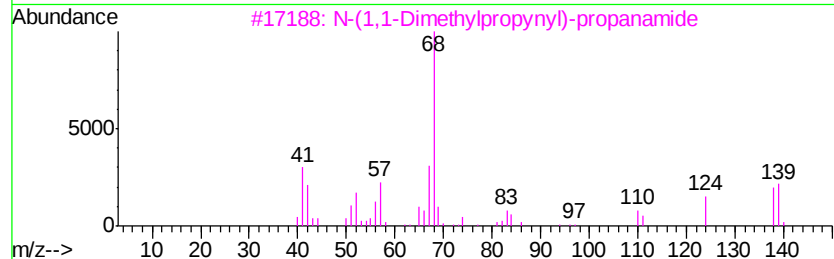
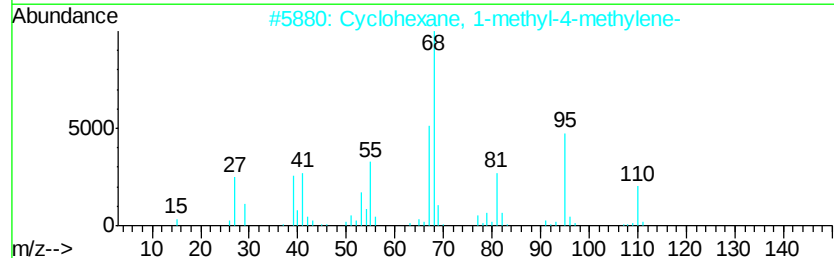
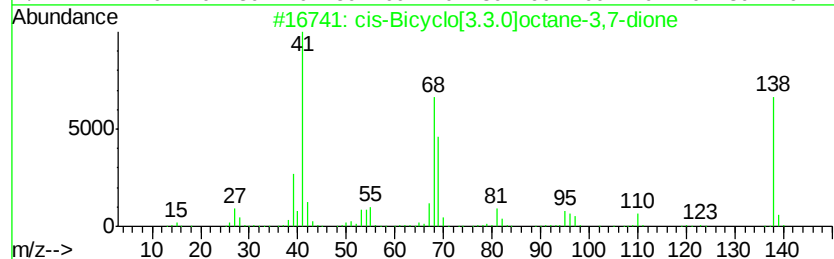
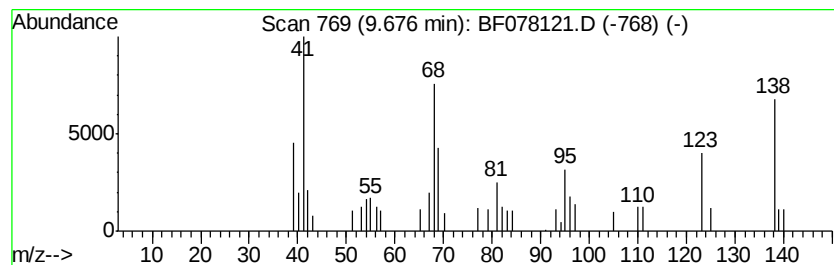
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 unknown9.68 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.68	3.95 ng	68061	Naphthalene-d8	8.62

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	cis-Bicyclo[3.3.0]octane-3,7-dione	138	C8H10O2	051716-63-3	49
2		Cyclohexane, 1-methyl-4-methylene-	110	C8H14	002808-80-2	35
3		N-(1,1-Dimethylpropynyl)-propana...	139	C8H13NO	1000110-54-3	35
4		Benzene, 1,2-dimethoxy-	138	C8H10O2	000091-16-7	35
5		Benzene, (ethylthio)-	138	C8H10S	000622-38-8	27



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF033115\
Data File : BF078121.D
Acq On : 31 Mar 2015 18:37
Operator : TP/IZ
Sample : G1700-01
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
S8-0424

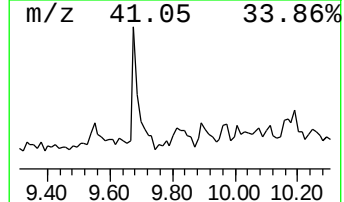
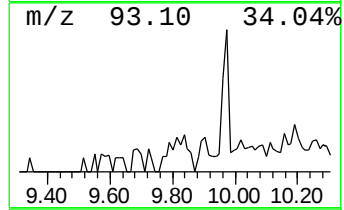
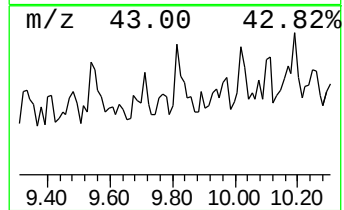
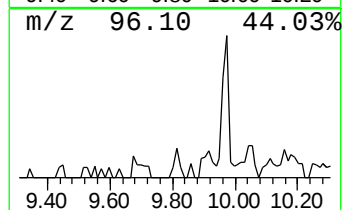
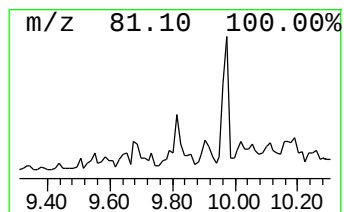
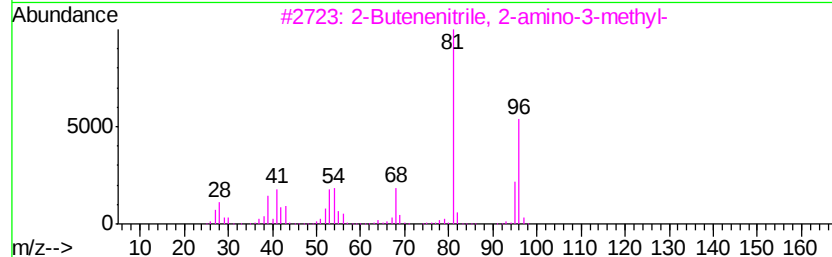
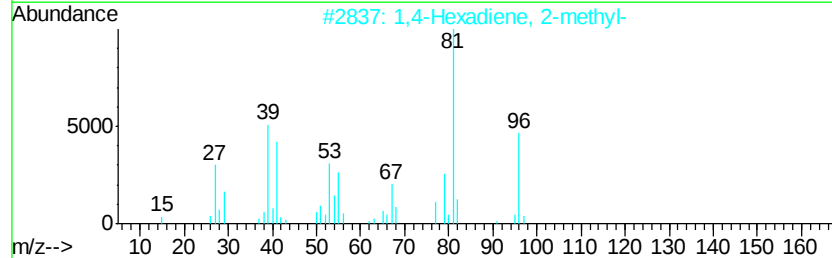
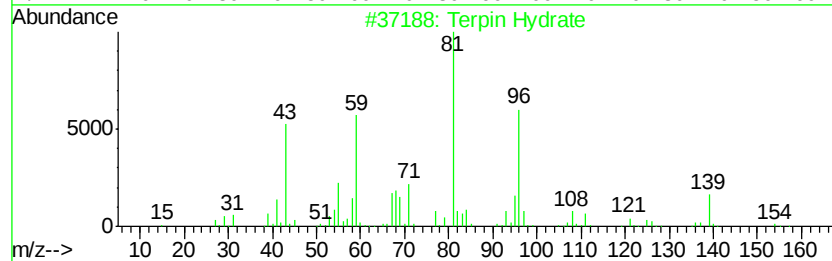
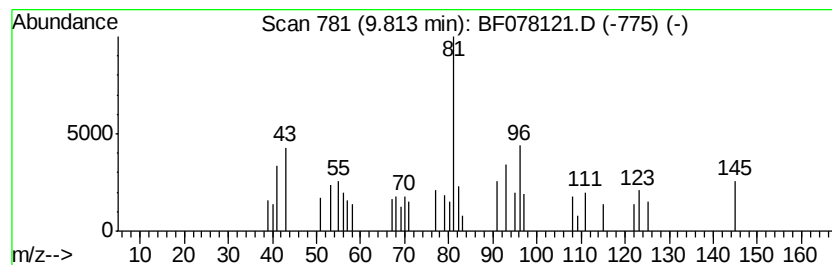
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 unknown9.81 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.81	2.67 ng	56454	Acenaphthene-d10	10.78

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Terpin Hydrate	172	C10H20O2	002451-01-6	43
2			1,4-Hexadiene, 2-methyl-	96	C7H12	001119-14-8	43
3			2-Butenenitrile, 2-amino-3-methyl-	96	C5H8N2	1000196-28-0	43
4			1,4-Hexadiene, 4-methyl-	96	C7H12	001116-90-1	43
5			3-Octyne, 2,2-dimethyl-	138	C10H18	019482-57-6	40



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF033115\
Data File : BF078121.D
Acq On : 31 Mar 2015 18:37
Operator : TP/IZ
Sample : G1700-01
Misc :
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ClientSampleId :
S8-0424

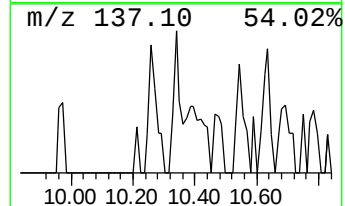
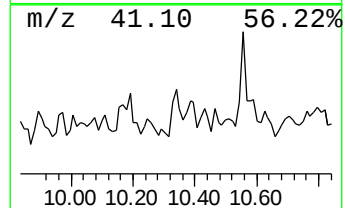
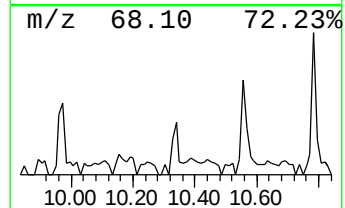
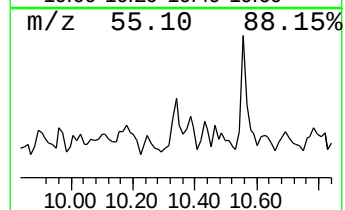
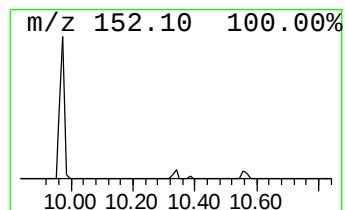
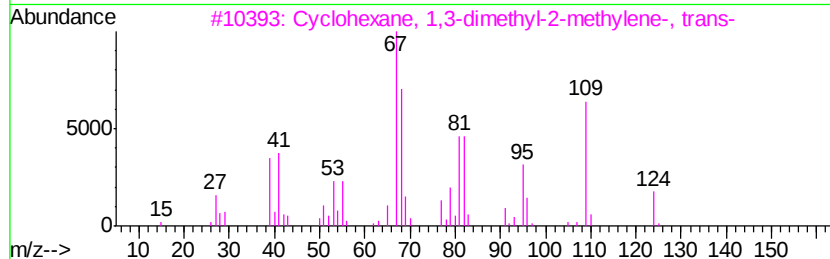
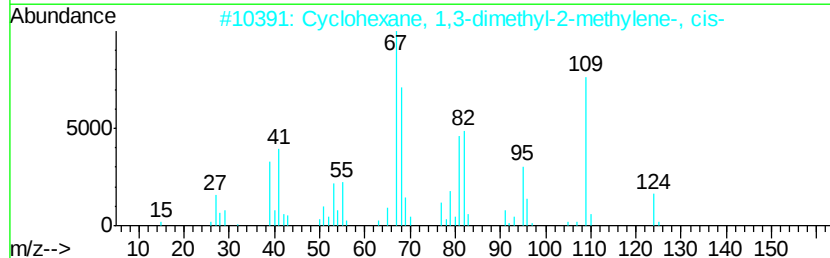
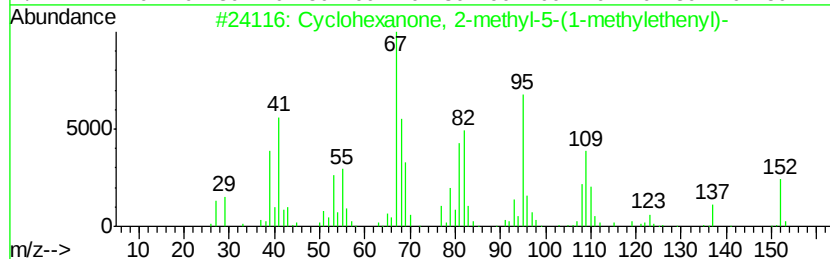
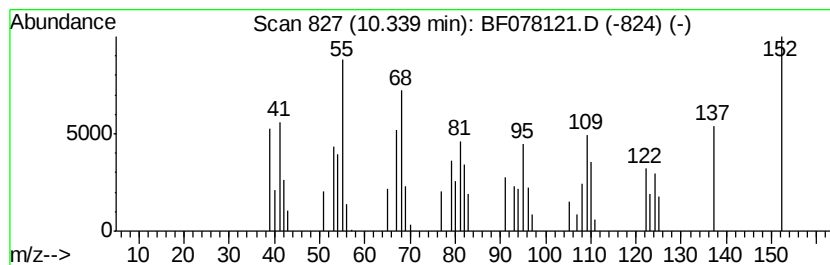
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 unknown10.34 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.34	2.01 ng	42395	Acenaphthene-d10	10.78

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexanone, 2-methyl-5-(1-met...	152	C10H16O	007764-50-3	49
2		Cyclohexane, 1,3-dimethyl-2-meth...	124	C9H16	019781-47-6	46
3		Cyclohexane, 1,3-dimethyl-2-meth...	124	C9H16	020348-74-7	45
4		3-Undecyne	152	C11H20	060212-30-8	43
5		Cyclopentanol, 1-(1-methylene-2-...	138	C9H14O	078158-11-9	38



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF031115\
Data File : BF078121.D
Acq On : 31 Mar 2015 18:37
Operator : TP/IZ
Sample : G1700-01
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
S8-0424

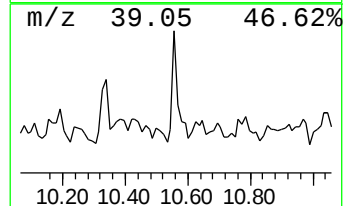
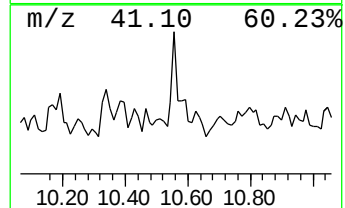
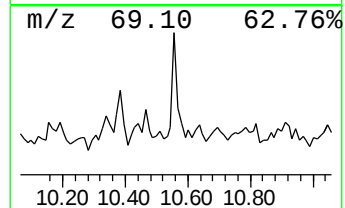
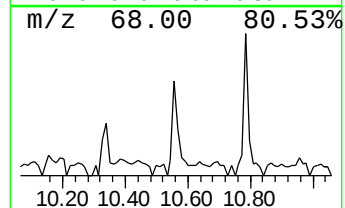
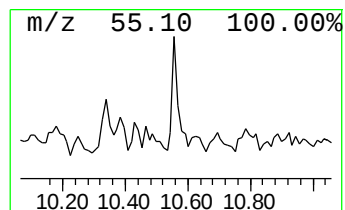
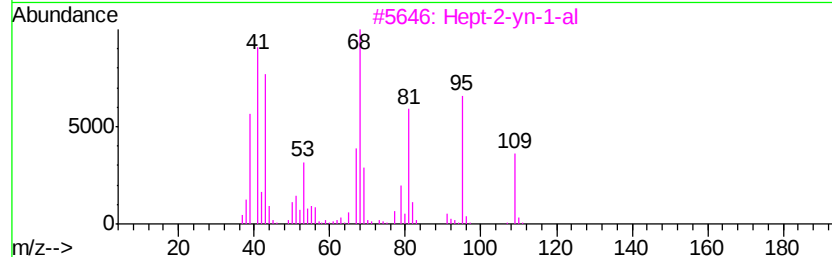
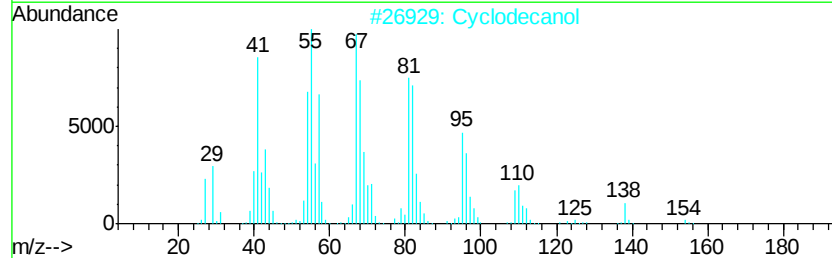
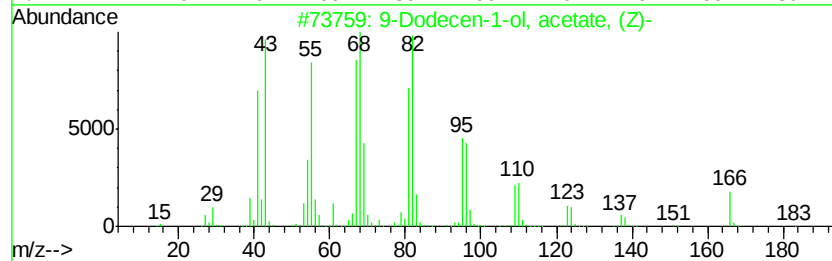
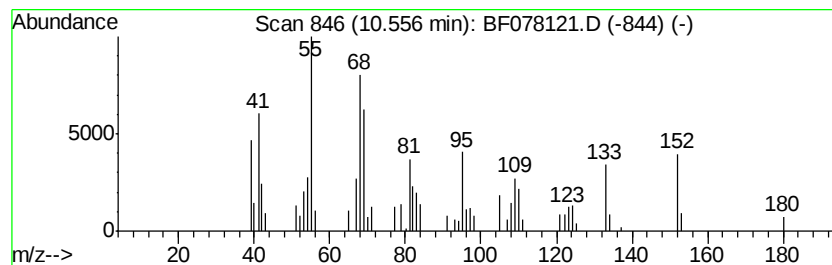
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 9-Dodecen-1-ol, acetate, (Z)- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.56	2.52 ng	53196	Acenaphthene-d10	10.78

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			9-Dodecen-1-ol, acetate, (Z)-	226	C14H26O2	016974-11-1	50
2			Cyclodecanol	156	C10H20O	001502-05-2	47
3			Hept-2-yn-1-al	110	C7H10O	001846-67-9	43
4			Bicyclo[3.3.1]nonane-2,8-dione	152	C9H12O2	160651-01-4	43
5			Z-1,6-Tridecadiene	180	C13H24	1000130-98-3	35



Data Path : Z:\HPCHEM1\BNA F\DATA\BF031115\
Data File : BF078121.D
Acq On : 31 Mar 2015 18:37
Operator : TP/IZ
Sample : G1700-01
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
S8-0424

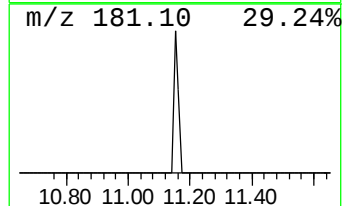
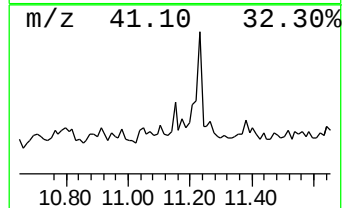
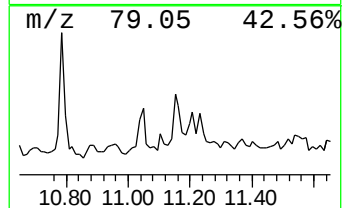
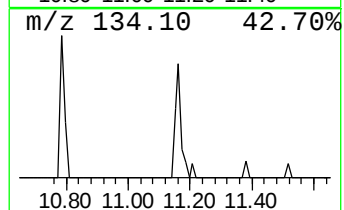
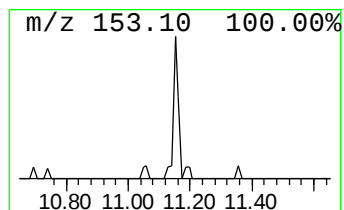
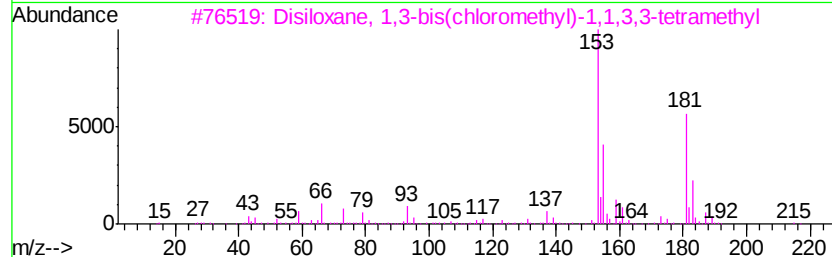
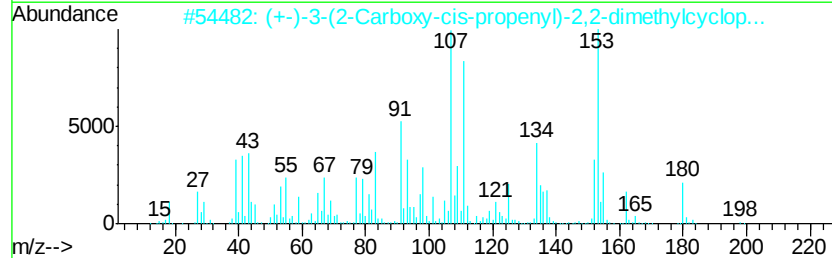
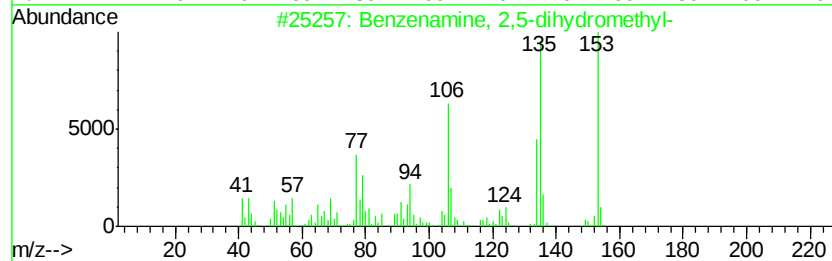
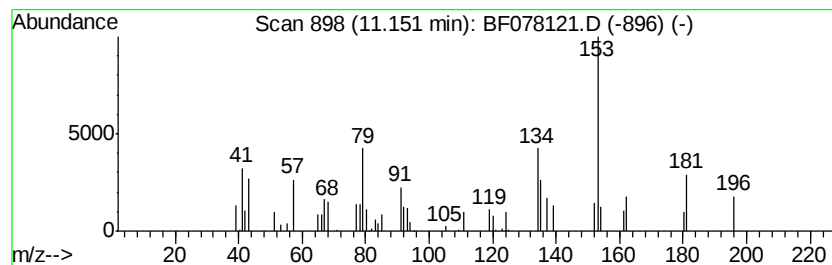
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 unknown11.15 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.15	3.71 ng	78369	Acenaphthene-d10	10.78

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzenamine, 2,5-dihydromethyl-	153	C8H11NO2	133242-65-6	47
2		(+)-3-(2-Carboxy-cis-propenyl)-...	198	C10H14O4	040253-02-9	38
3		Disiloxane, 1,3-bis(chloromethyl-...	230	C6H16Cl2OSi2	002362-10-9	35
4		(+)-3-(2-Carboxy-cis-propenyl)-2...	198	C10H14O4	056390-95-5	28
5		Benzenamine, 3,4-dimethoxy-	153	C8H11NO2	006315-89-5	25



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF031115\
Data File : BF078121.D
Acq On : 31 Mar 2015 18:37
Operator : TP/IZ
Sample : G1700-01
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
S8-0424

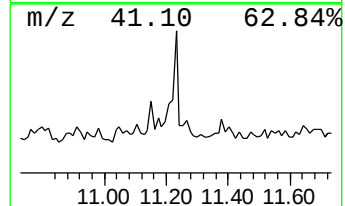
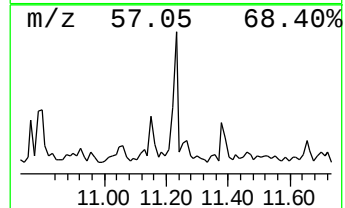
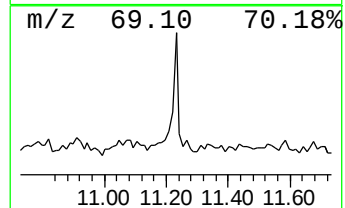
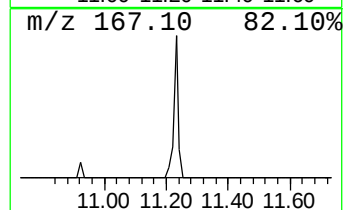
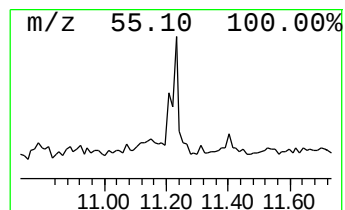
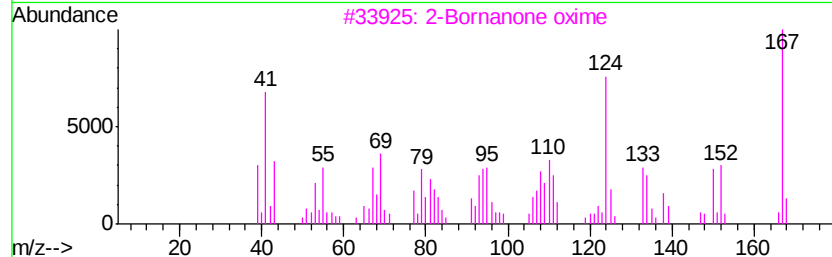
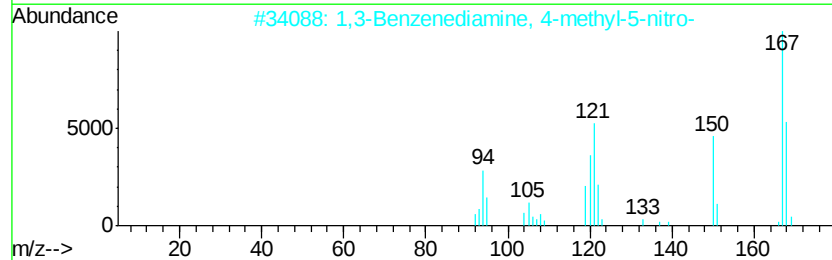
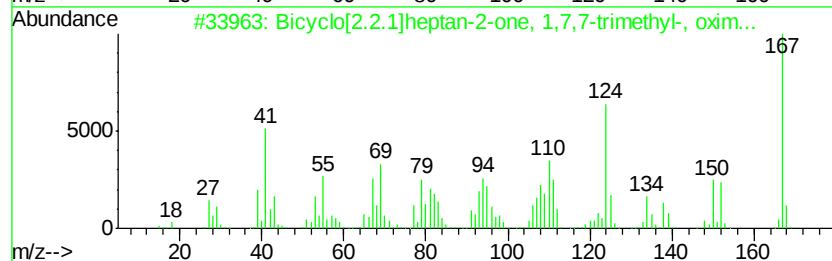
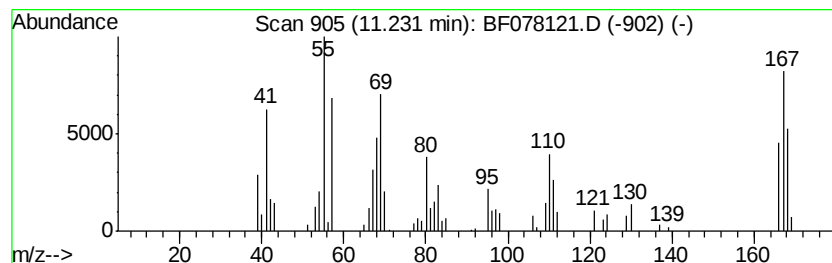
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 unknown11.23 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.23	5.07 ng	107054	Acenaphthene-d10	10.78

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Bicyclo[2.2.1]heptan-2-one, 1,7,...	167	C10H17NO	018674-50-5	43
2			1,3-Benzenediamine, 4-methyl-5-n...	167	C7H9N3O2	006629-29-4	38
3			2-Bornanone oxime	167	C10H17NO	013559-66-5	25
4			2,7(1H,3H)-Naphthalenedione, hex...	166	C10H14O2	046048-64-0	22
5			Chloromethyl 6-chloroheptanoate	212	C8H14Cl2O2	080418-62-8	22



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF033115\
Data File : BF078121.D
Acq On : 31 Mar 2015 18:37
Operator : TP/IZ
Sample : G1700-01
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031115.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...	4.77	3.4	ng	40157	1	7.05	237311	20.0
unknown6.76	6.76	83.9	ng	995079	1	7.05	237311	20.0
unknown8.11	8.11	2.9	ng	49830	2	8.62	344960	20.0
unknown9.68	9.68	4.0	ng	68061	2	8.62	344960	20.0
unknown9.81	9.81	2.7	ng	56454	3	10.78	422325	20.0
unknown10.34	10.34	2.0	ng	42395	3	10.78	422325	20.0
9-Dodecen-1-ol, a...	10.56	2.5	ng	53196	3	10.78	422325	20.0
unknown11.15	11.15	3.7	ng	78369	3	10.78	422325	20.0
unknown11.23	11.23	5.1	ng	107054	3	10.78	422325	20.0