

Data Path : Z:\HPCHEM1\BNA F\DATA\BF032515\
 Data File : BF077968.D
 Acq On : 25 Mar 2015 23:13
 Operator : TP/IZ
 Sample : G1643-06
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW-7-3-23-15

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	1.356	38	41	50	rVB	188672	335116	2.23%	0.471%
2	1.516	53	55	58	rVV	233154	310759	2.07%	0.437%
3	1.630	62	65	74	rVV	243146	398452	2.66%	0.560%
4	1.744	74	75	103	rVB	644886	1130983	7.54%	1.589%
5	5.379	390	393	400	rVB	358791	450874	3.01%	0.633%
6	6.693	506	508	516	rVB	164239	277226	1.85%	0.389%
7	6.807	516	518	521	rBV	740029	928693	6.19%	1.304%
8	6.876	522	524	526	rBV	525562	718779	4.79%	1.010%
9	6.922	526	528	530	rVB	292282	336427	2.24%	0.473%
10	7.059	536	540	541	rBV2	2996712	4176033	27.85%	5.866%
11	7.093	541	543	545	rVB	286728	257945	1.72%	0.362%
12	7.173	547	550	552	rBV	1262244	1081485	7.21%	1.519%
13	7.288	557	560	562	rBV	677467	911004	6.08%	1.280%
14	7.790	602	604	607	rVB	838059	761632	5.08%	1.070%
15	8.602	671	675	679	rBV2	220869	378686	2.53%	0.532%
16	8.671	679	681	684	rVB	415587	365857	2.44%	0.514%
17	9.265	730	733	736	rVB	410379	702672	4.69%	0.987%
18	9.345	736	740	741	rBV	3494034	6354444	42.37%	8.926%
19	9.379	741	743	745	rVV	3904171	8824643	58.85%	12.395%
20	9.413	745	746	748	rVV	3403088	2508334	16.73%	3.523%
21	9.505	749	754	755	rBV	5899495	14995941	100.00%	21.064%
22	9.539	755	757	760	rVB2	6397751	9869073	65.81%	13.863%
23	9.642	762	766	768	rBV	4906766	7134890	47.58%	10.022%
24	10.031	797	800	802	rVB	1330813	1481489	9.88%	2.081%
25	10.442	833	836	839	rVB	138813	237077	1.58%	0.333%
26	10.591	845	849	851	rBV	817774	1419538	9.47%	1.994%
27	10.671	854	856	858	rBV	322943	317317	2.12%	0.446%
28	10.842	869	871	873	rVB	425364	420994	2.81%	0.591%
29	11.814	954	956	959	rBV	865166	882478	5.88%	1.240%
30	12.671	1029	1031	1034	rBV	462772	451580	3.01%	0.634%
31	14.671	1203	1206	1208	rBV	1639370	1669464	11.13%	2.345%
32	15.951	1315	1318	1320	rBV	481331	545268	3.64%	0.766%
33	17.620	1461	1464	1467	rBV	487965	557321	3.72%	0.783%

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Instrument :
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ClientSampleId :
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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_F\METHODS\8270-BF031115.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

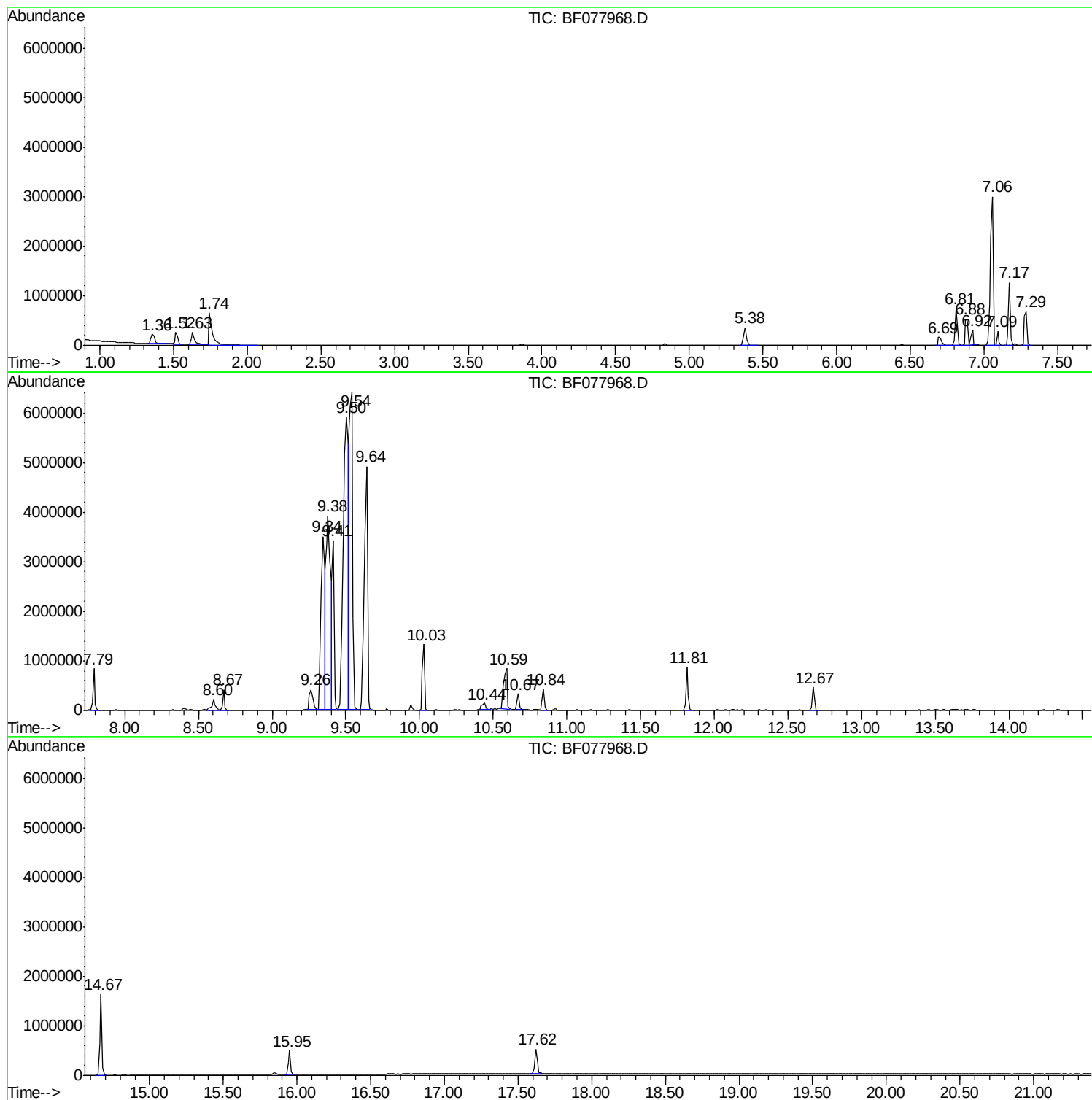
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF031115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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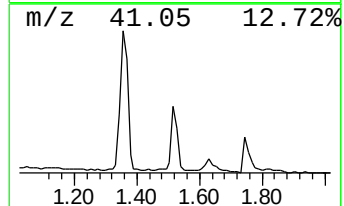
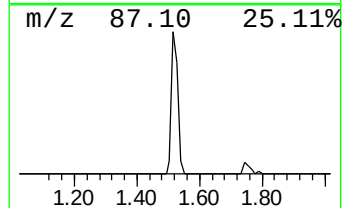
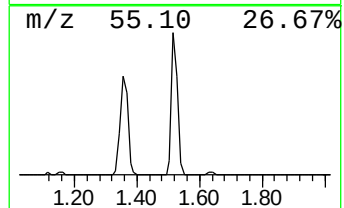
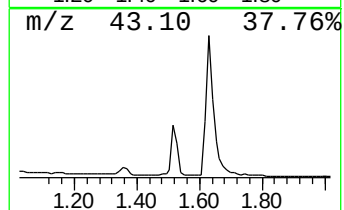
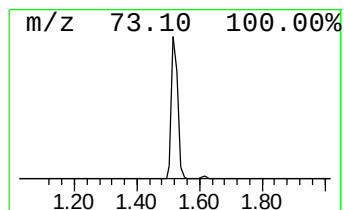
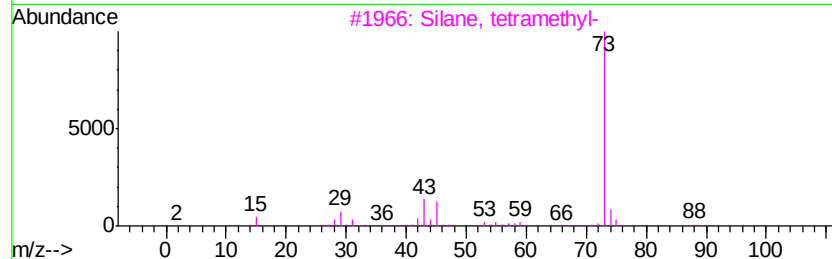
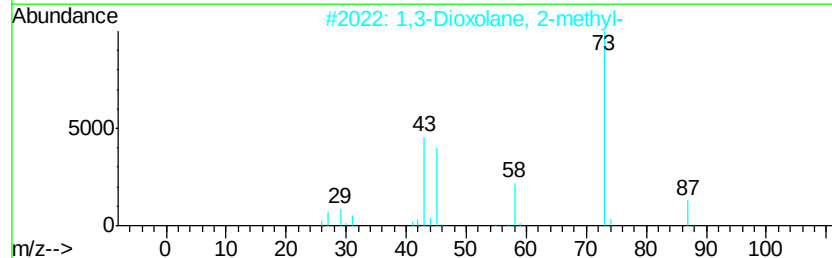
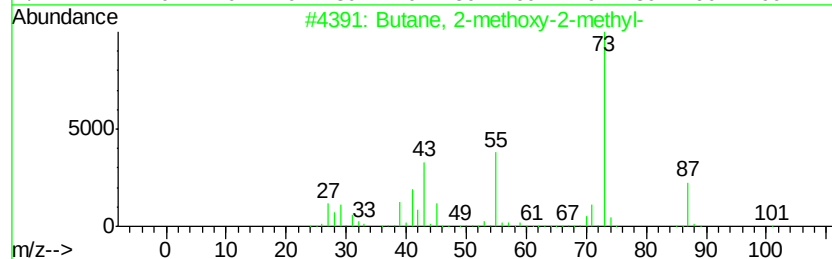
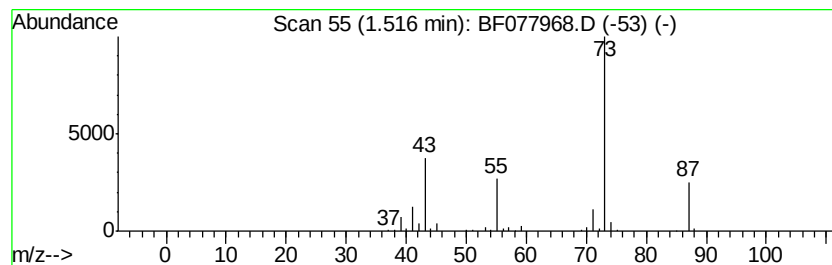
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Butane, 2-methoxy-2-methyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.52	24.10 ng	310759	1,4-Dichlorobenzene-d4	7.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2		1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	37
3		Silane, tetramethyl-	88	C4H12Si	000075-76-3	25
4		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	23
5		Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	10



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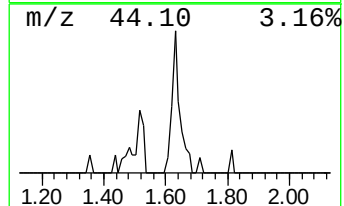
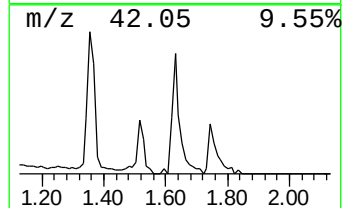
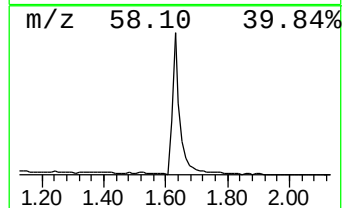
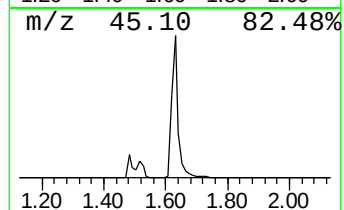
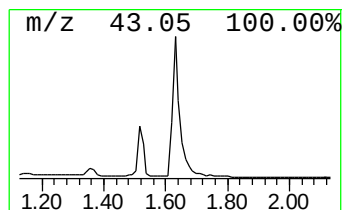
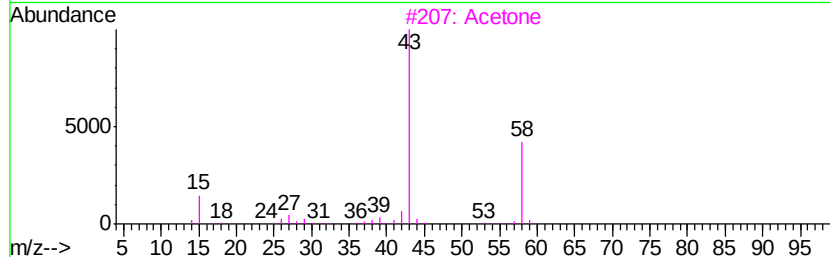
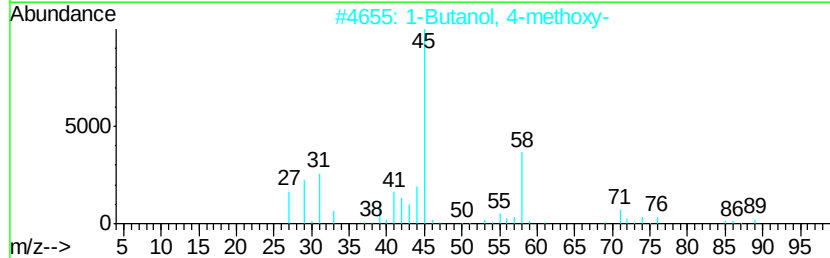
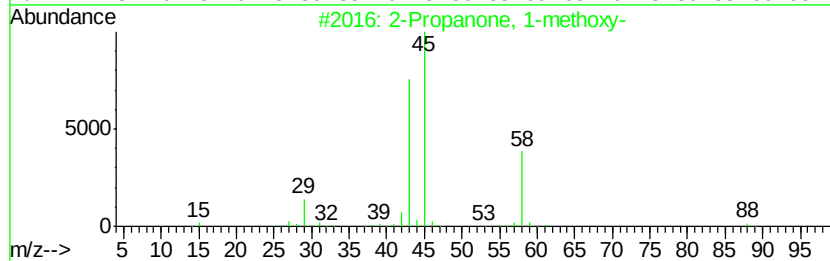
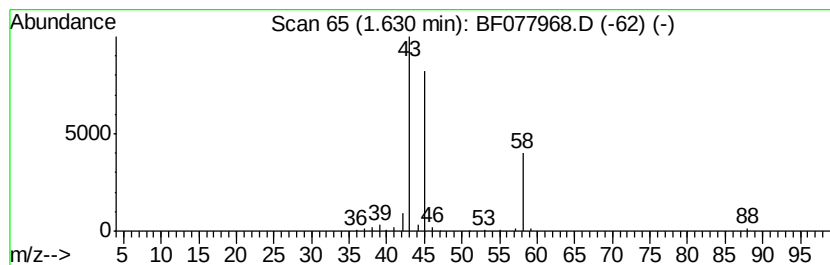
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 2-Propanone, 1-methoxy- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.63	30.89 ng	398452	1,4-Dichlorobenzene-d4	7.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Propanone, 1-methoxy-	88	C4H8O2	005878-19-3	86
2		1-Butanol, 4-methoxy-	104	C5H12O2	000111-32-0	50
3		Acetone	58	C3H6O	000067-64-1	38
4		Butane, 1-methoxy-	88	C5H12O	000628-28-4	9
5		Ethanol, 2-methoxy-, acetate	118	C5H10O3	000110-49-6	9



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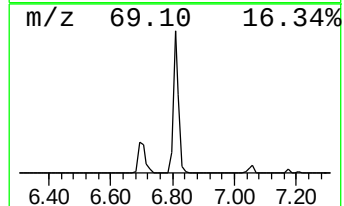
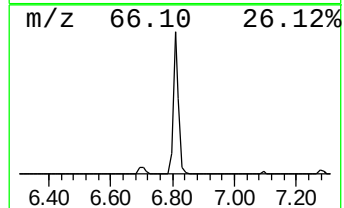
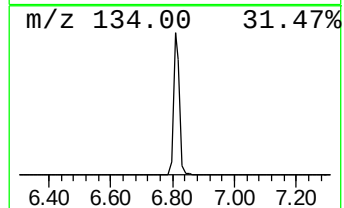
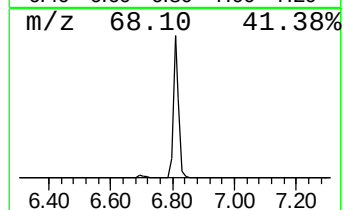
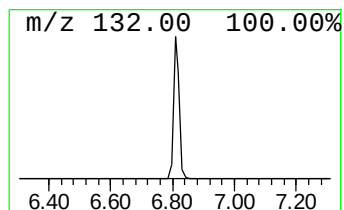
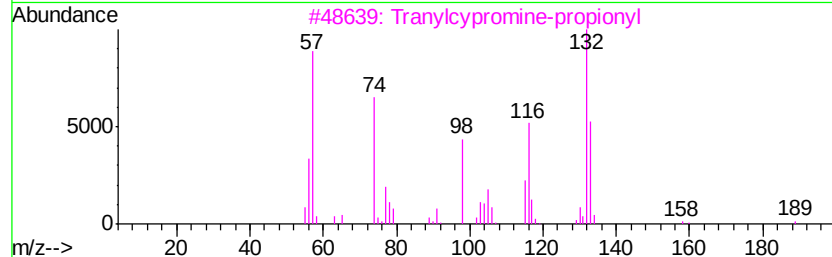
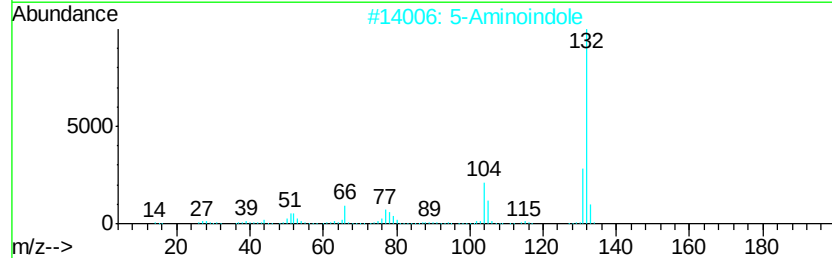
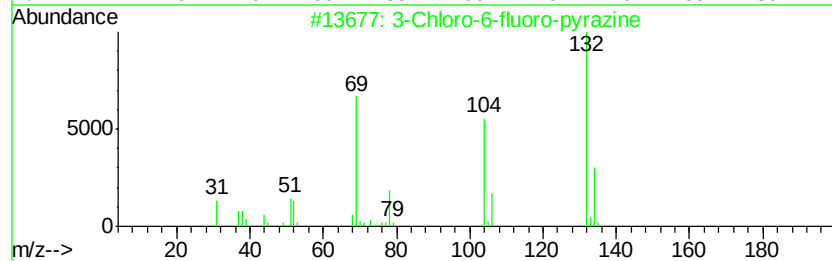
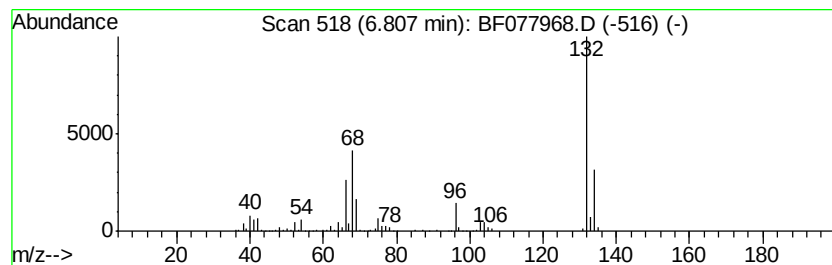
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 unknown6.81 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.81	72.01 ng	928693	1,4-Dichlorobenzene-d4	7.09

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2	5-Aminoindole	132	C8H8N2	005192-03-0	12
3	Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
4	4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9
5	1-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-59-9	9



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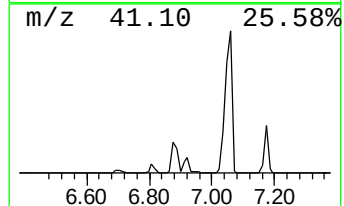
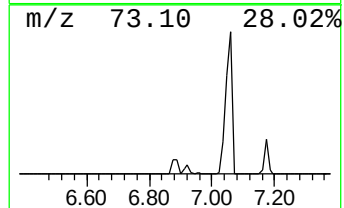
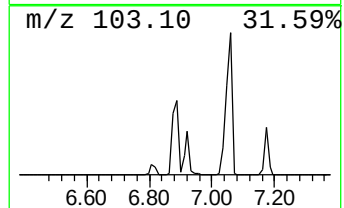
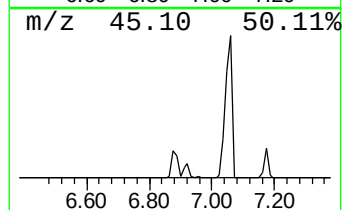
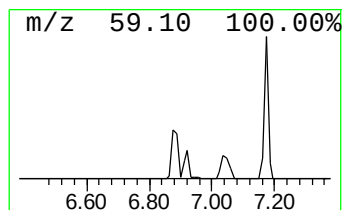
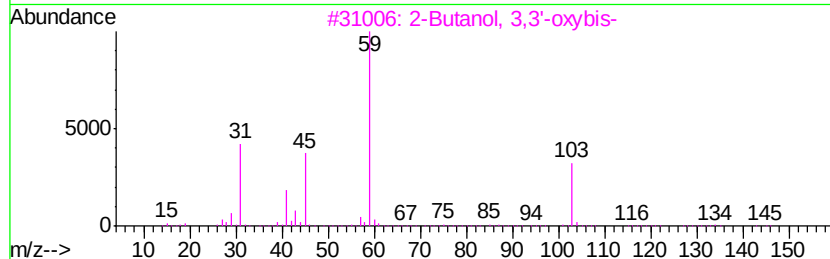
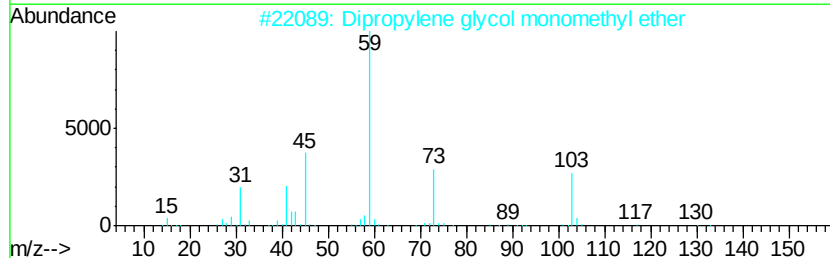
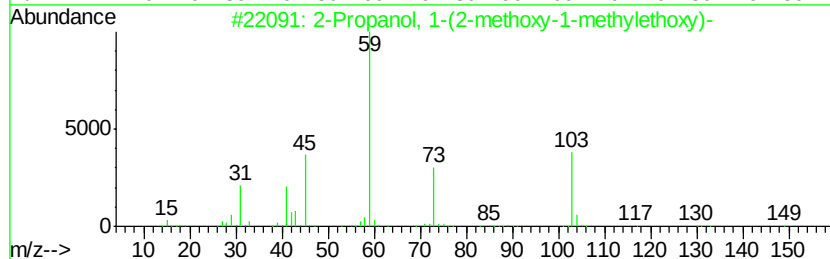
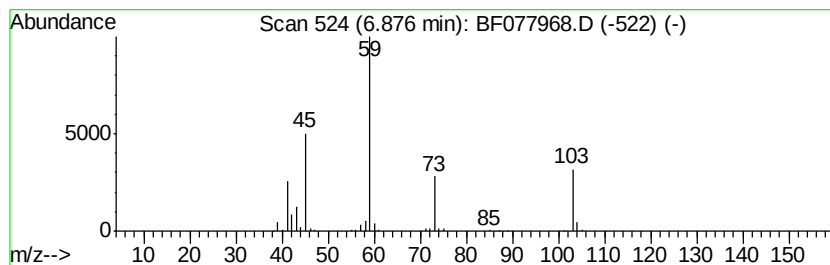
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 2-Propanol, 1-(2-methoxy-1-... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.88	55.73 ng	718779	1,4-Dichlorobenzene-d4	7.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Propanol, 1-(2-methoxy-1-methy...	148	C7H16O3	020324-32-7	80
2		Dipropylene glycol monomethyl ether	148	C7H16O3	034590-94-8	72
3		2-Butanol, 3,3'-oxybis-	162	C8H18O3	054305-61-2	64
4		1-Propanol, 2,2'-oxybis-	134	C6H14O3	000108-61-2	53
5		1-Propanol, 2-(2-hydroxypropoxy)-	134	C6H14O3	000106-62-7	45



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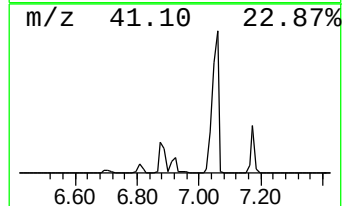
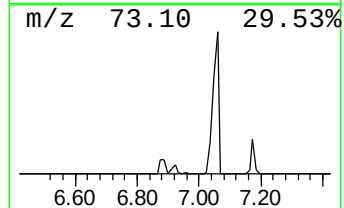
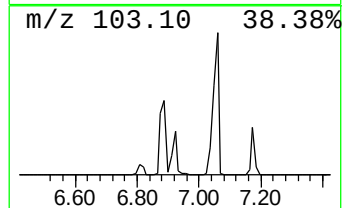
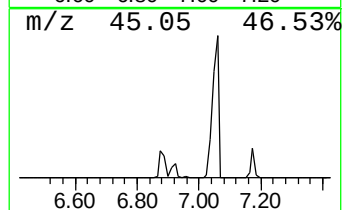
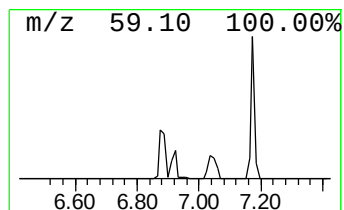
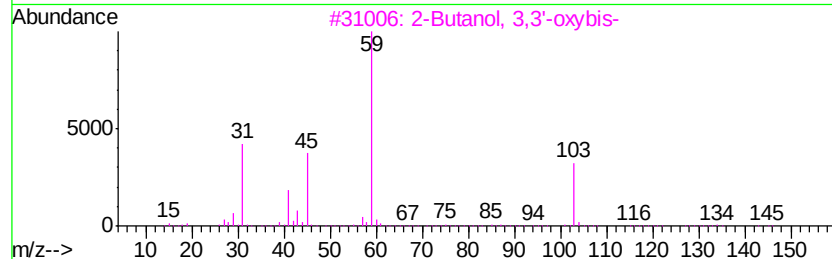
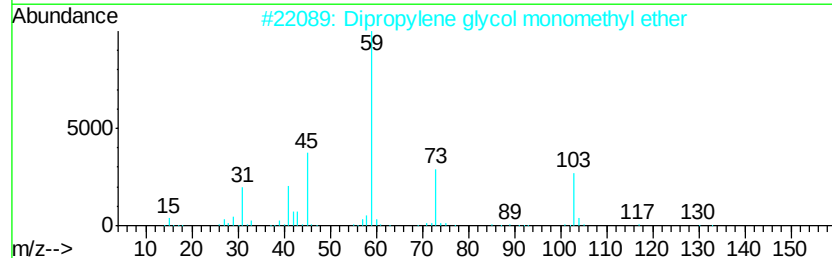
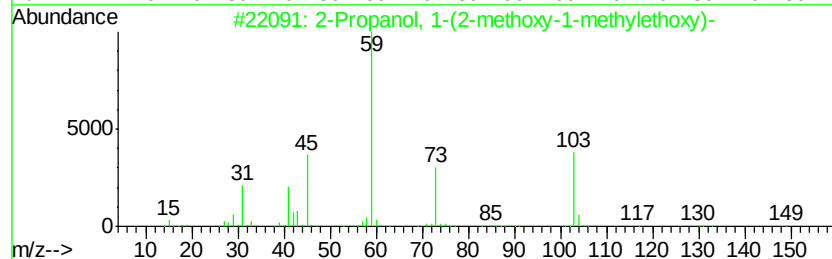
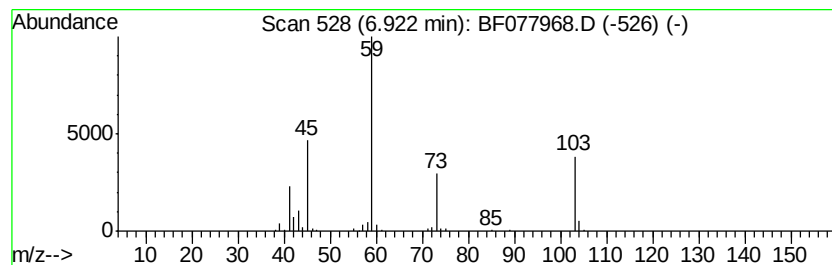
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Dipropylene glycol monometh... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.92	26.09 ng	336427	1,4-Dichlorobenzene-d4	7.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Propanol, 1-(2-methoxy-1-methy...	148	C7H16O3	020324-32-7	90
2		Dipropylene glycol monomethyl ether	148	C7H16O3	034590-94-8	72
3		2-Butanol, 3,3'-oxybis-	162	C8H18O3	054305-61-2	64
4		1-Propanol, 2-(2-hydroxypropoxy)-	134	C6H14O3	000106-62-7	45
5		1-Propanol, 2,2'-oxybis-	134	C6H14O3	000108-61-2	38



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MW-7-3-23-15

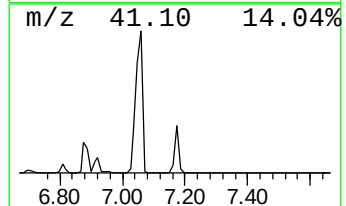
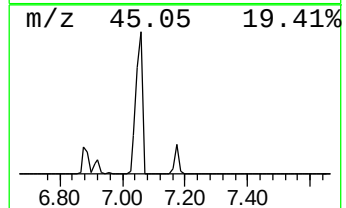
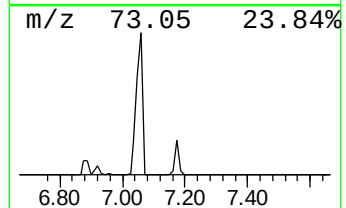
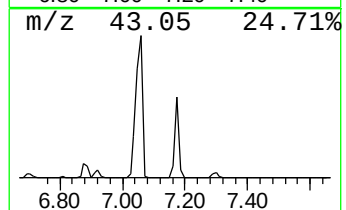
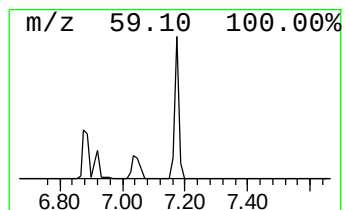
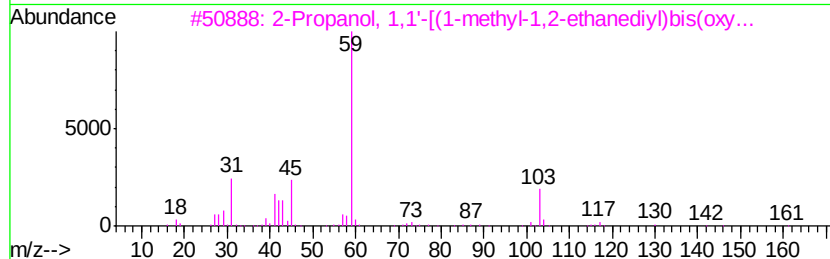
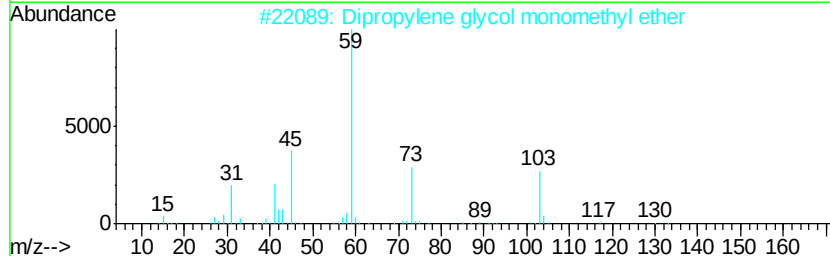
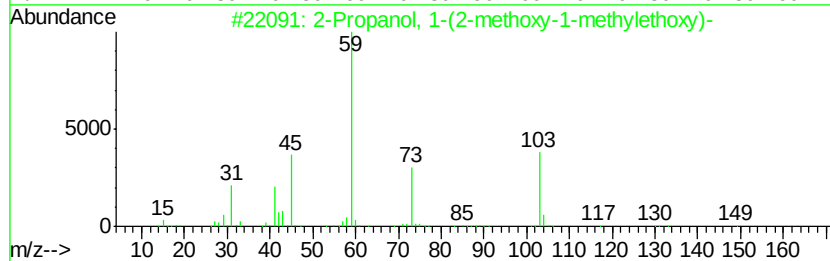
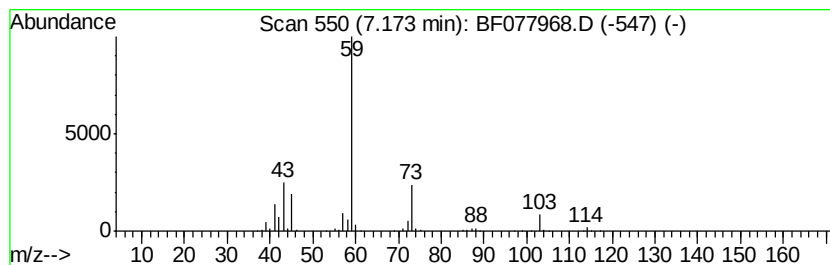
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 2-Propanol, 1,1'-[(1-methyl... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.17	83.85 ng	1081490	1,4-Dichlorobenzene-d4	7.09

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Propanol,	1-(2-methoxy-1-methy...	148	C7H16O3	020324-32-7	72
2	Dipropylene glycol	monomethyl ether	148	C7H16O3	034590-94-8	72
3	2-Propanol, 1,1'-[	(1-methyl-1,2-...	192	C9H20O4	001638-16-0	64
4	2-Butanol, 3-methoxy-		104	C5H12O2	053778-72-6	59
5	Butane, 2-methoxy-		88	C5H12O	006795-87-5	59



Data Path : Z:\HPCHEM1\BNA F\DATA\BF032515\
Data File : BF077968.D
Acq On : 25 Mar 2015 23:13
Operator : TP/IZ
Sample : G1643-06
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MW-7-3-23-15

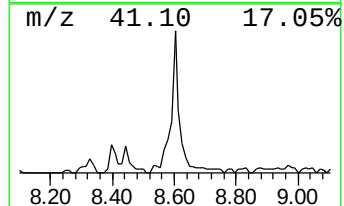
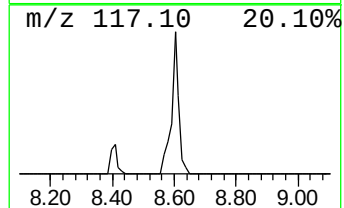
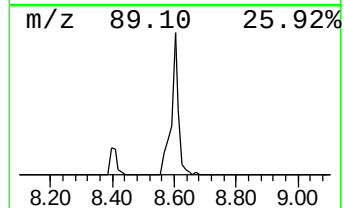
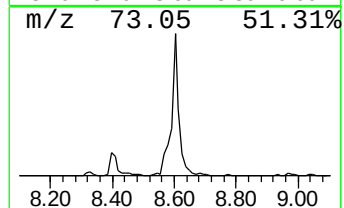
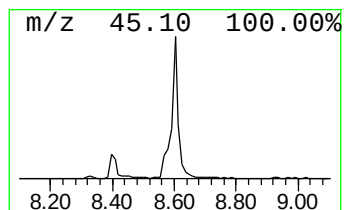
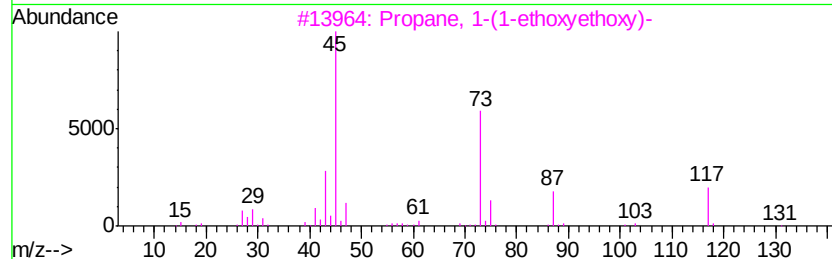
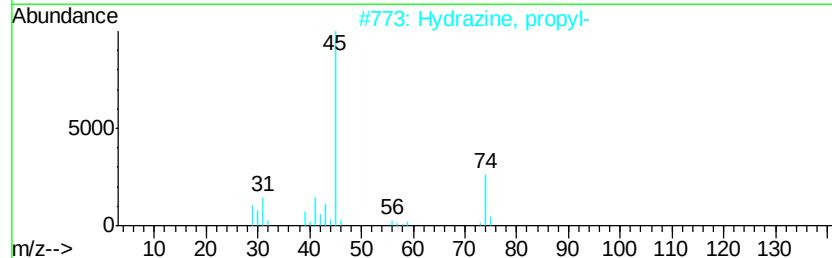
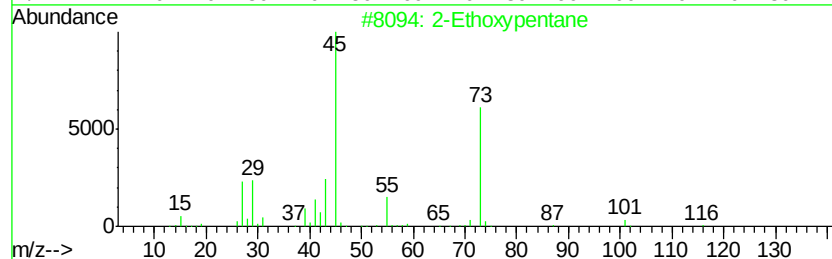
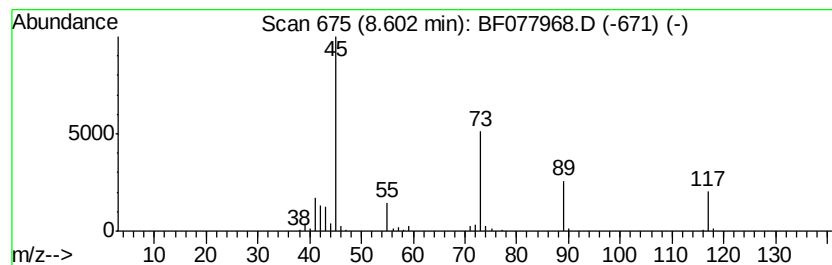
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 unknown8.60 Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.60	20.70 ng	378686	Naphthalene-d8	8.67

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Ethoxypentane	116	C7H16O	001817-89-6	40
2		Hydrazine, propyl-	74	C3H10N2	005039-61-2	38
3		Propane, 1-(1-ethoxvethoxy)-	132	C7H16O2	020680-10-8	38
4		Hexane, 1-methoxy-	116	C7H16O	004747-07-3	28
5		Pentane, 1-methoxy-	102	C6H14O	000628-80-8	28



Data Path : Z:\HPCHEM1\BNA F\DATA\BF032515\
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Operator : TP/IZ
Sample : G1643-06
Misc :
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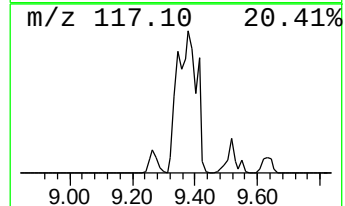
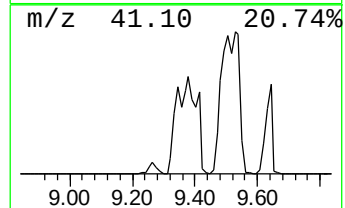
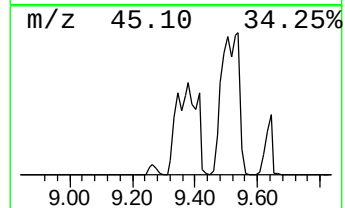
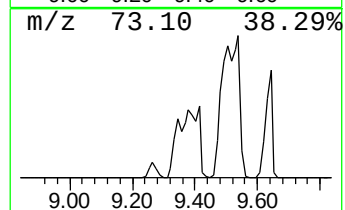
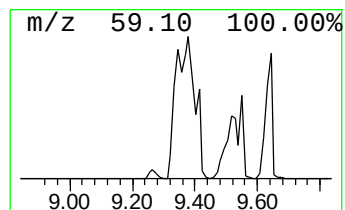
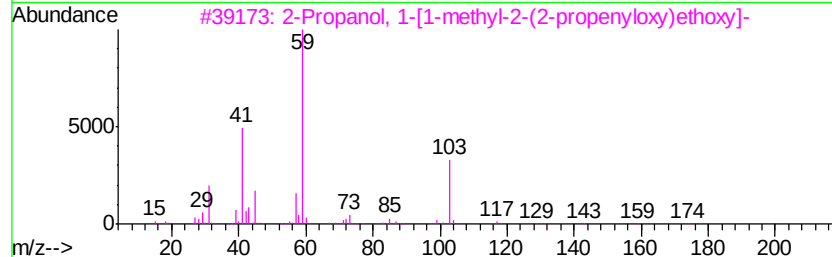
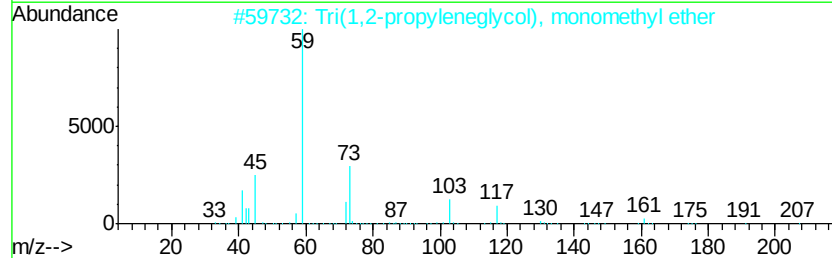
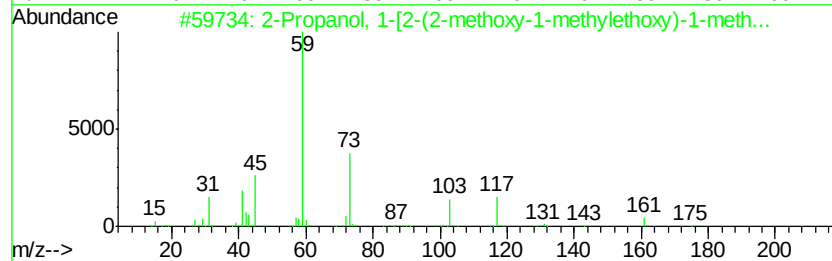
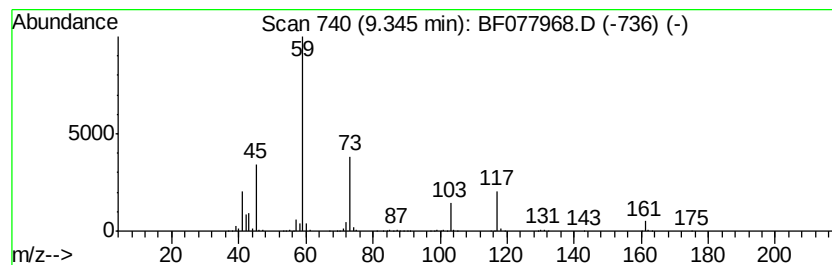
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 2-Propanol, 1-[2-(2-methoxy... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.34	347.37 ng	6354440	Napthalene-d8	8.67
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	2-Propanol, 1-[2-(2-methoxy-1-me...	206 C10H22O4	020324-33-8	90
2	Tri(1,2-propyleneglycol), monome...	206 C10H22O4	1000262-26-6	64
3	2-Propanol, 1-[1-methyl-2-(2-pro...	174 C9H18O3	055956-25-7	53
4	1-Propanol, 2-(2-hydroxypropoxy)-	134 C6H14O3	000106-62-7	50
5	2-Propanol, 1-(2-methoxy-1-methy...	148 C7H16O3	020324-32-7	50



Data Path : Z:\HPCHEM1\BNA F\DATA\BF032515\
Data File : BF077968.D
Acq On : 25 Mar 2015 23:13
Operator : TP/IZ
Sample : G1643-06
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
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ClientSampleId :
MW-7-3-23-15

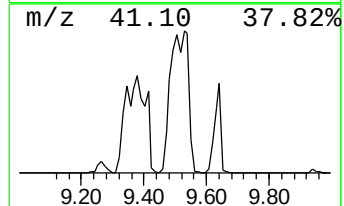
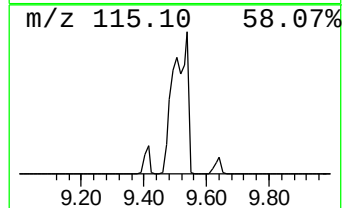
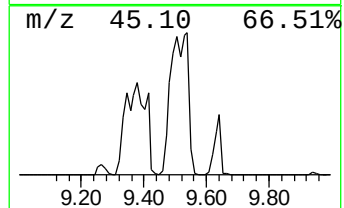
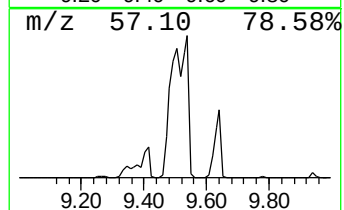
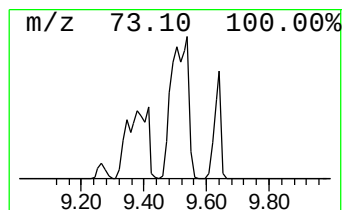
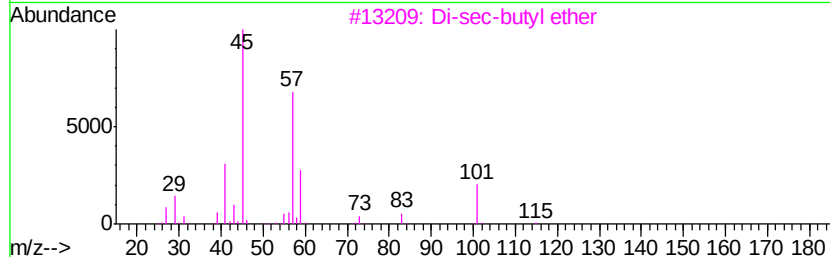
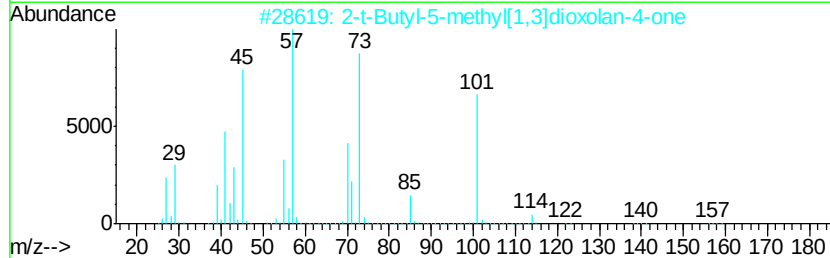
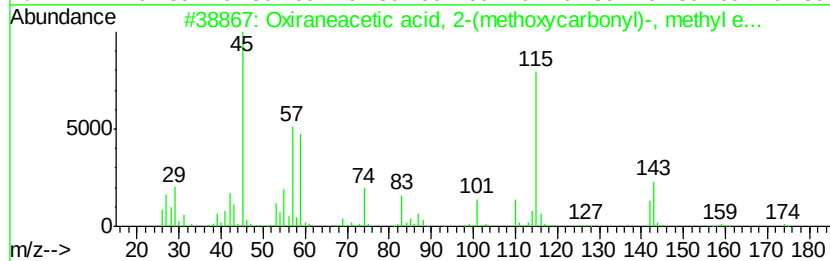
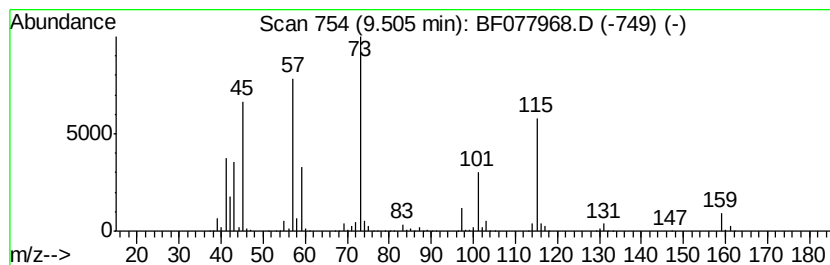
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 unknown9.50 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.50	819.77 ng	14995900	Napthalene-d8	8.67

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Oxiraneacetic acid, 2-(methoxycar...	174	C7H10O5	052336-07-9	38
2		2-t-Butyl-5-methyl[1,3]dioxolan-...	158	C8H14O3	130930-47-1	38
3		Di-sec-butyl ether	130	C8H18O	006863-58-7	37
4		Butane, 2-ethoxy-	102	C6H14O	002679-87-0	35
5		Butane, 1-(1-methylpropoxy)-	130	C8H18O	000999-65-5	25



Data Path : Z:\HPCHEM1\BNA F\DATA\BF032515\
Data File : BF077968.D
Acq On : 25 Mar 2015 23:13
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Sample : G1643-06
Misc :
ALS Vial : 17 Sample Multiplier: 1

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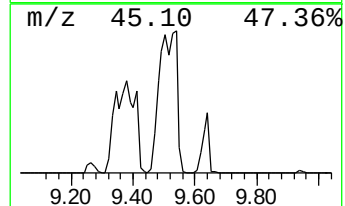
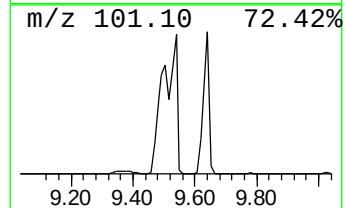
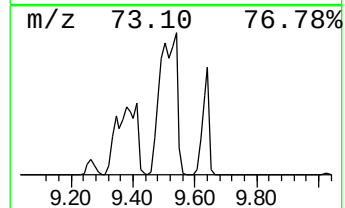
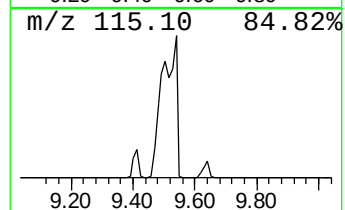
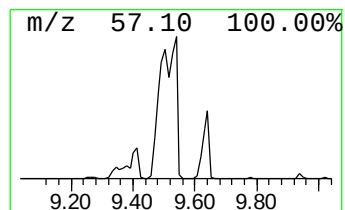
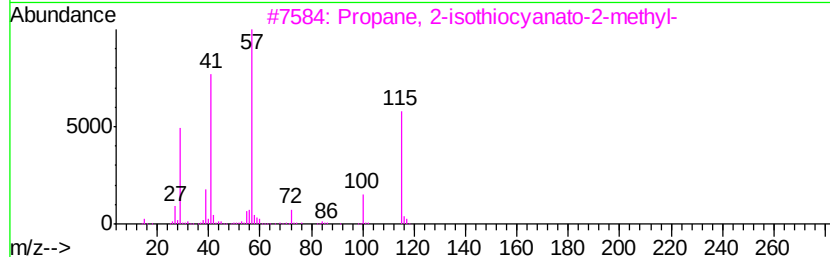
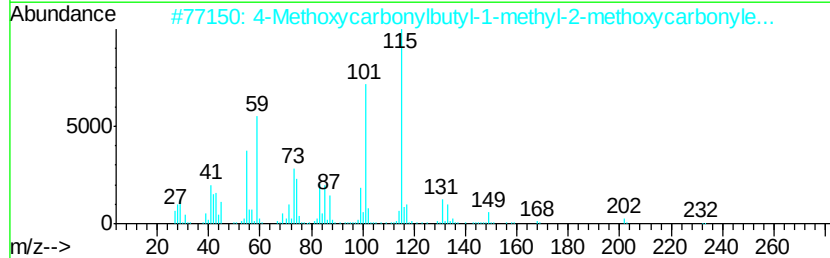
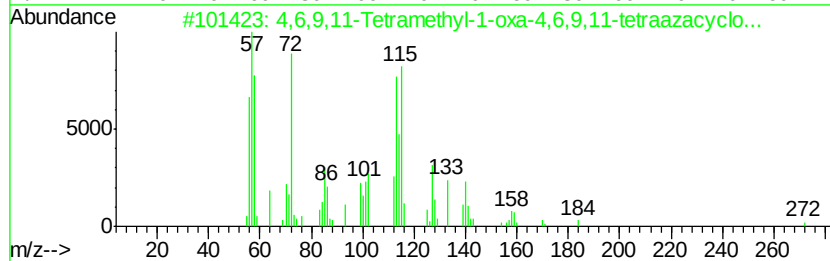
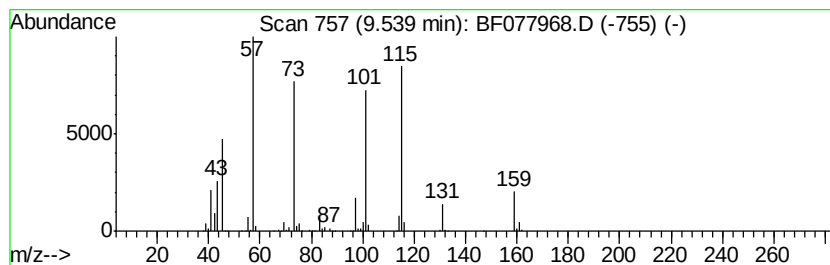
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 unknown9.54 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.54	539.50 ng	9869070	Napthalene-d8	8.67

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4,6,9,11-Tetramethyl-1-oxa-4,6,9...	272	C12H24N4O3	1000139-36-0	38
2		4-Methoxycarbonylbutyl-1-methyl-...	232	C11H20O5	1000140-30-2	28
3		Propane, 2-isothiocyanato-2-methyl-	115	C5H9NS	000590-42-1	27
4		2-Oxopentanedioic acid	146	C5H6O5	000328-50-7	12
5		1,3,5-Cyclohexanetriol, (1.alpha...	132	C6H12O3	050409-12-6	11



Data Path : Z:\HPCHEM1\BNA F\DATA\BF032515\
Data File : BF077968.D
Acq On : 25 Mar 2015 23:13
Operator : TP/IZ
Sample : G1643-06
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
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ClientSampleId :
MW-7-3-23-15

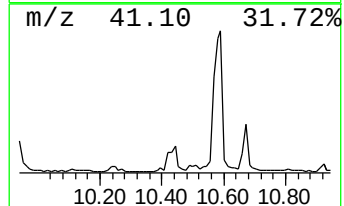
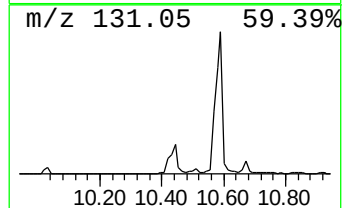
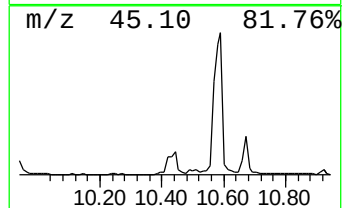
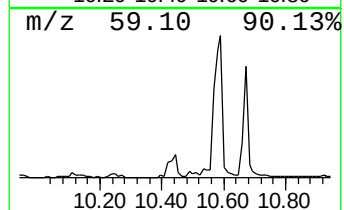
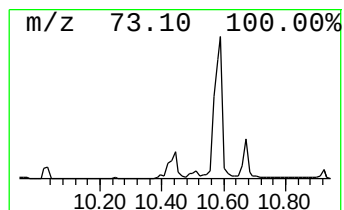
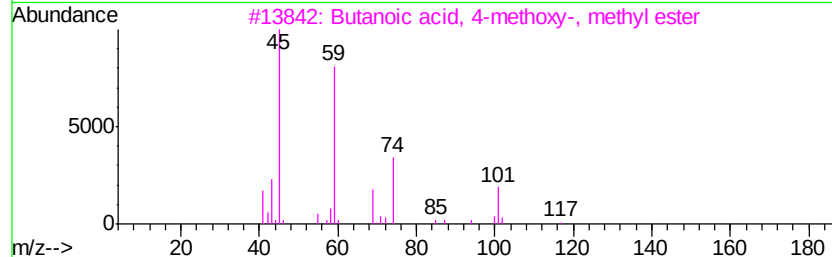
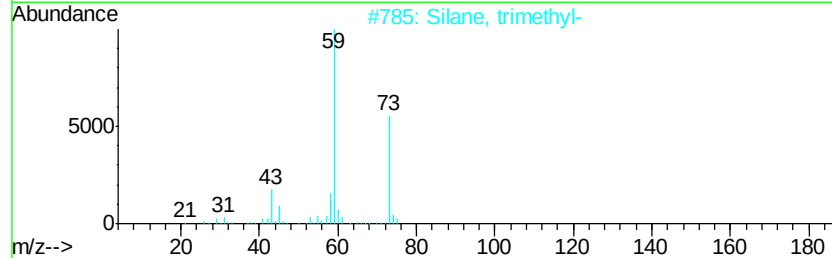
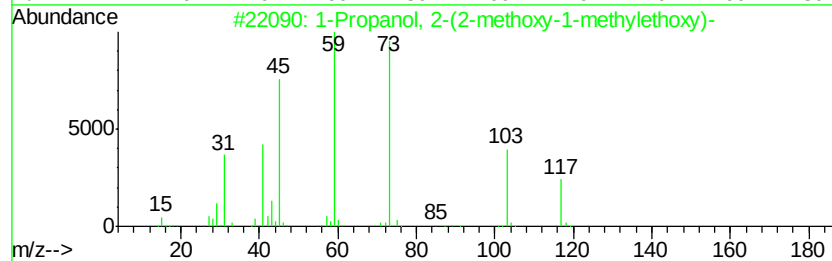
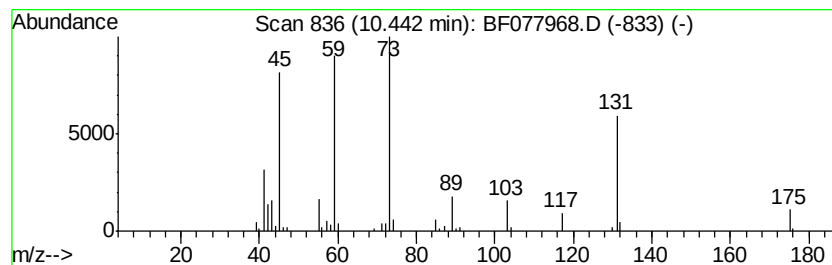
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 unknown10.44 Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.44	11.26 ng	237077	Acenaphthene-d10	10.84

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Propanol, 2-(2-methoxy-1-methy...	148	C7H16O3	055956-21-3	43
2			Silane, trimethyl-	74	C3H10Si	000993-07-7	40
3			Butanoic acid, 4-methoxy-, methy...	132	C6H12O3	029006-01-7	38
4			Ethane, 1-ethoxy-1-methoxy-	104	C5H12O2	010471-14-4	37
5			2-Propanol, 1-[2-(2-methoxy-1-me...	206	C10H22O4	020324-33-8	35



Data Path : Z:\HPCHEM1\BNA F\DATA\BF032515\
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Misc :
ALS Vial : 17 Sample Multiplier: 1

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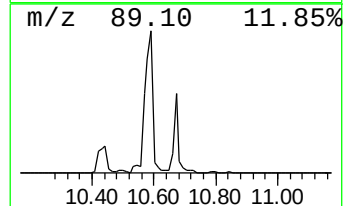
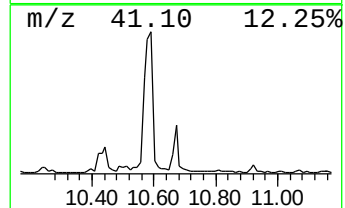
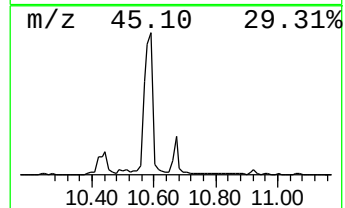
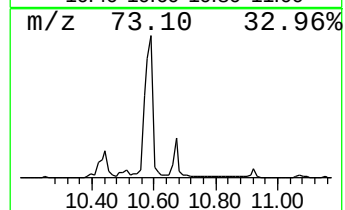
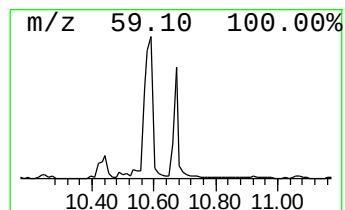
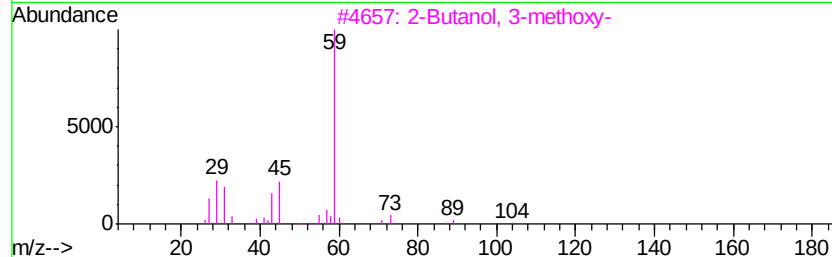
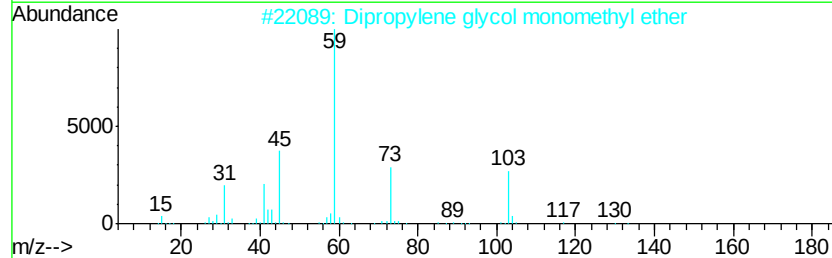
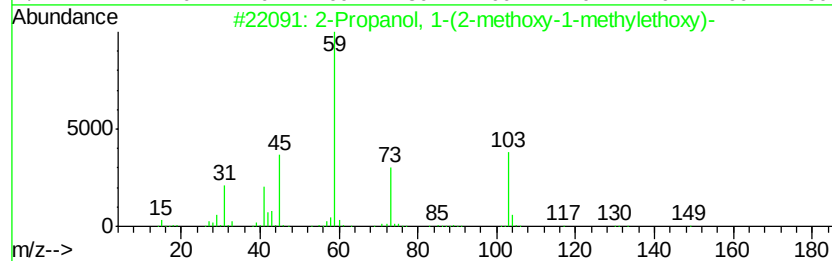
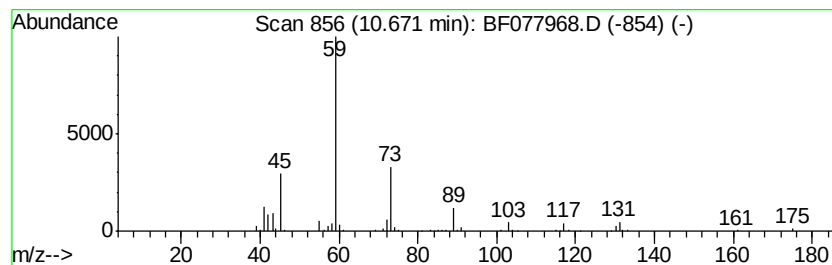
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 19 2-Butanol, 3-methoxy- Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.67	15.07 ng	317317	Acenaphthene-d10	10.84

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Propanol, 1-(2-methoxy-1-methyl-...	148	C7H16O3	020324-32-7	64
2		Dipropylene glycol monomethyl ether	148	C7H16O3	034590-94-8	64
3		2-Butanol, 3-methoxy-	104	C5H12O2	053778-72-6	53
4		2-Propanol, 1-methoxy-2-methyl-	104	C5H12O2	003587-64-2	50
5		2-Propanol, 1-[2-(2-methoxy-1-me...	206	C10H22O4	020324-33-8	50



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032515\
 Data File : BF077968.D
 Acq On : 25 Mar 2015 23:13
 Operator : TP/IZ
 Sample : G1643-06
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW-7-3-23-15

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
Butane, 2-methoxy...	1.52	24.1	ng	310759	1	7.09	257945	20.0
2-Propanone, 1-me...	1.63	30.9	ng	398452	1	7.09	257945	20.0
unknown6.81	6.81	72.0	ng	928693	1	7.09	257945	20.0
2-Propanol, 1-(2-...	6.88	55.7	ng	718779	1	7.09	257945	20.0
Dipropylene glyco...	6.92	26.1	ng	336427	1	7.09	257945	20.0
2-Propanol, 1,1'-...	7.17	83.8	ng	1081490	1	7.09	257945	20.0
unknown8.60	8.60	20.7	ng	378686	2	8.67	365857	20.0
2-Propanol, 1-[2-...	9.34	347.4	ng	6354440	2	8.67	365857	20.0
unknown9.50	9.50	819.8	ng	14995900	2	8.67	365857	20.0
unknown9.54	9.54	539.5	ng	9869070	2	8.67	365857	20.0
unknown10.44	10.44	11.3	ng	237077	3	10.84	420994	20.0
2-Butanol, 3-meth...	10.67	15.1	ng	317317	3	10.84	420994	20.0