

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF021216\
 Data File : BF085037.D
 Acq On : 12 Feb 2016 3:16
 Operator : SJ/IZ
 Sample : H1410-01
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SUP-B004(16-18)

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-BF020116.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	1.735	11	13	14	rBV	1208373	1620426	17.00%	2.978%
2	1.850	21	23	25	rVB	7003173	9532452	100.00%	17.518%
3	2.261	57	59	66	rVB	117575	158995	1.67%	0.292%
4	4.753	274	277	288	rVB	847771	1347101	14.13%	2.476%
5	5.141	308	311	314	rBV	178845	338989	3.56%	0.623%
6	5.210	314	317	319	rBV	2853107	4055314	42.54%	7.452%
7	6.124	395	397	398	rBV	195211	154741	1.62%	0.284%
8	6.204	402	404	407	rVB	142289	119453	1.25%	0.220%
9	6.273	407	410	412	rBV	3550851	4410075	46.26%	8.104%
10	6.376	417	419	422	rBV	3446541	4003460	42.00%	7.357%
11	6.433	422	424	426	rBV	698066	585452	6.14%	1.076%
12	6.604	437	439	441	rBV	910321	863813	9.06%	1.587%
13	6.662	441	444	446	rVB	154578	172507	1.81%	0.317%
14	6.764	450	453	455	rBV	3714729	3279746	34.41%	6.027%
15	6.936	466	468	473	rVB	147709	138293	1.45%	0.254%
16	7.085	478	481	485	rVB	98940	161727	1.70%	0.297%
17	7.187	485	490	492	rBV	2198847	3281155	34.42%	6.030%
18	7.393	505	508	509	rBV	100786	129622	1.36%	0.238%
19	7.416	509	510	512	rVB	177848	130967	1.37%	0.241%
20	7.565	521	523	527	rVB	81386	129052	1.35%	0.237%
21	7.633	527	529	532	rBV3	235180	287016	3.01%	0.527%
22	7.896	549	552	556	rBV2	1413986	1695887	17.79%	3.117%
23	8.605	612	614	617	rBV	217035	196294	2.06%	0.361%
24	8.708	620	623	625	rVB	111090	116562	1.22%	0.214%
25	8.982	643	647	649	rBV	4814893	5413533	56.79%	9.948%
26	9.645	702	705	707	rBV	1464318	1326806	13.92%	2.438%
27	10.433	771	774	777	rBV	2958388	3084550	32.36%	5.668%
28	11.119	832	834	838	rBV	1245167	1264996	13.27%	2.325%
29	12.719	970	974	976	rBV	3293212	4801525	50.37%	8.824%
30	13.748	1062	1064	1067	rBV	928712	860451	9.03%	1.581%
31	15.108	1180	1183	1186	rBV	695619	754755	7.92%	1.387%

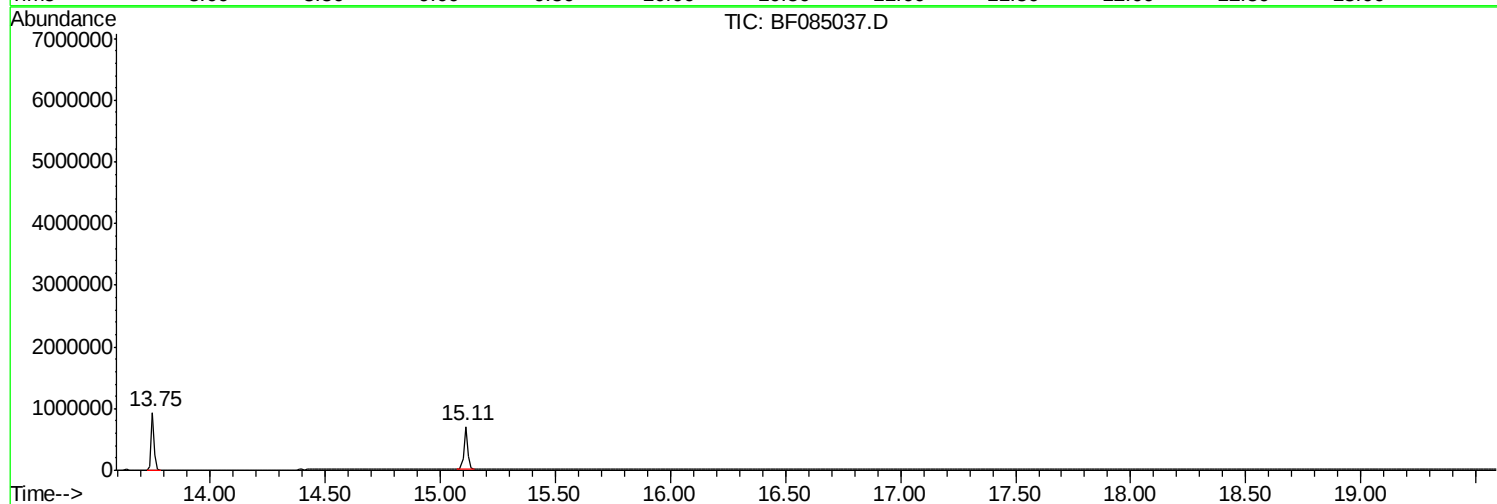
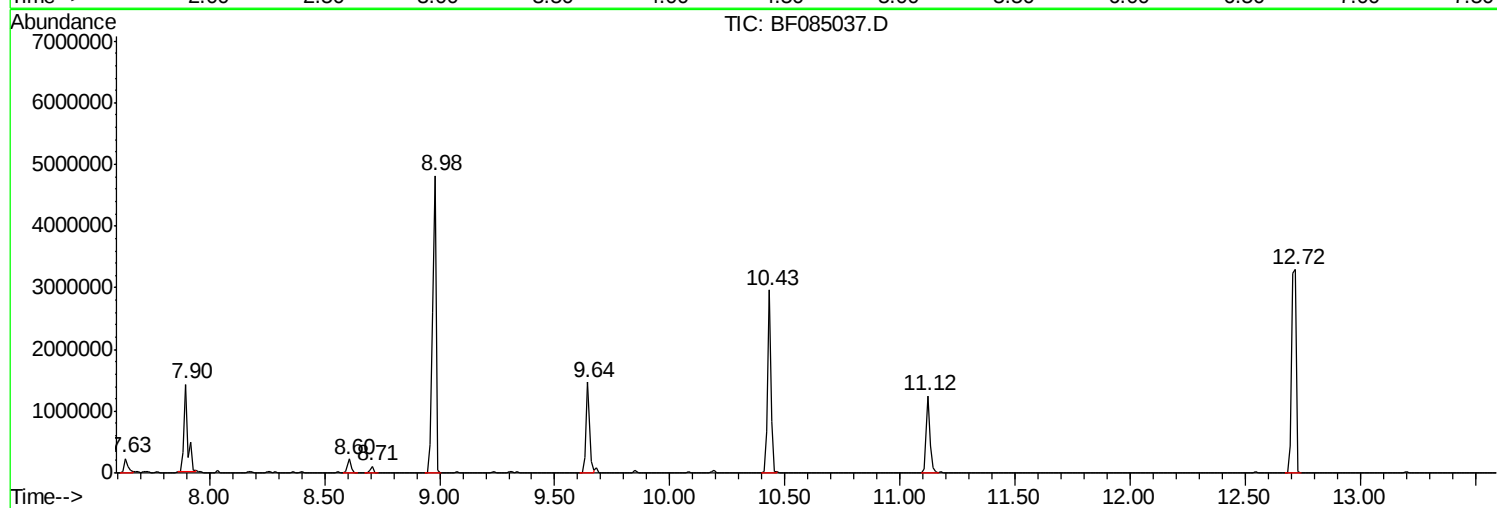
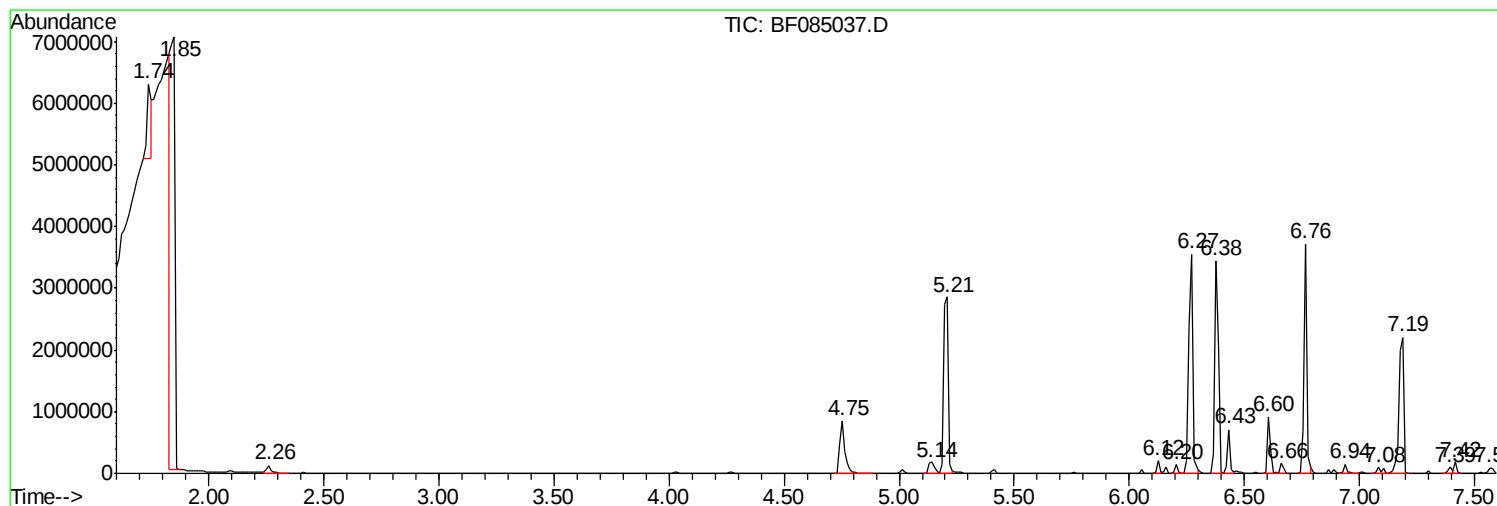
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Instrument :
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ClientSampleId :
SUP-B004(16-18)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF020116.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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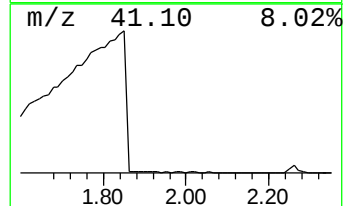
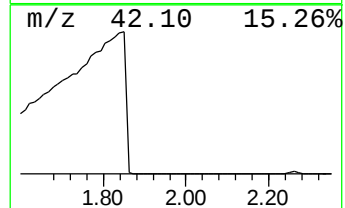
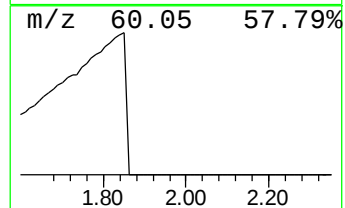
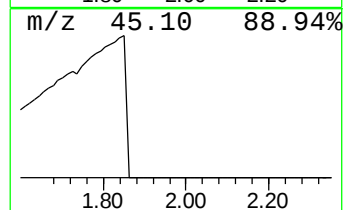
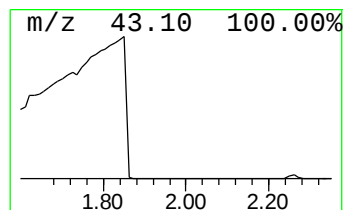
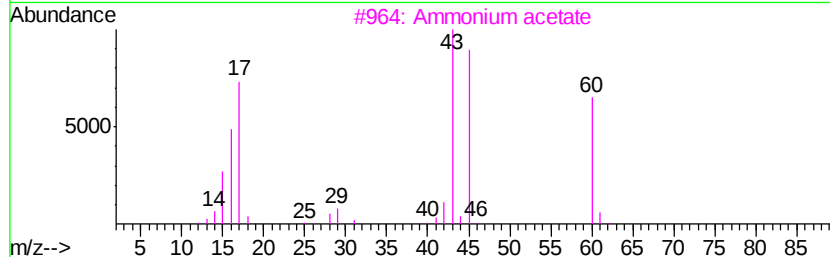
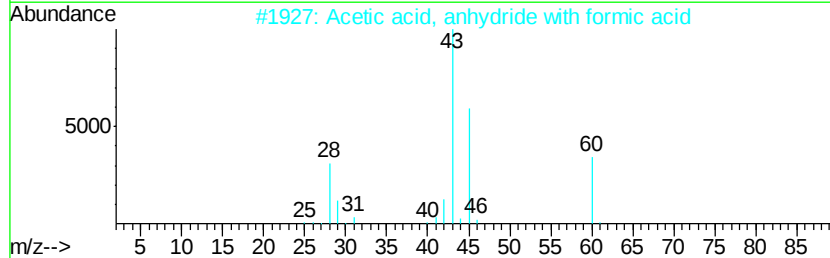
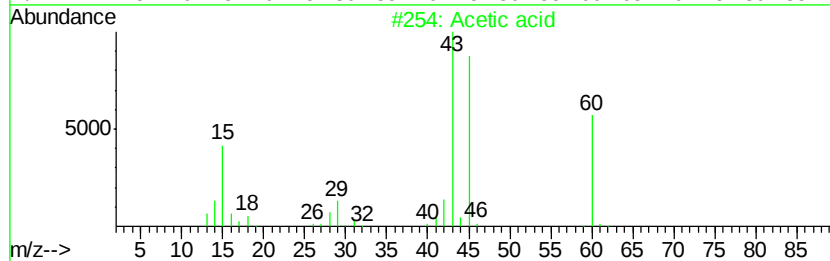
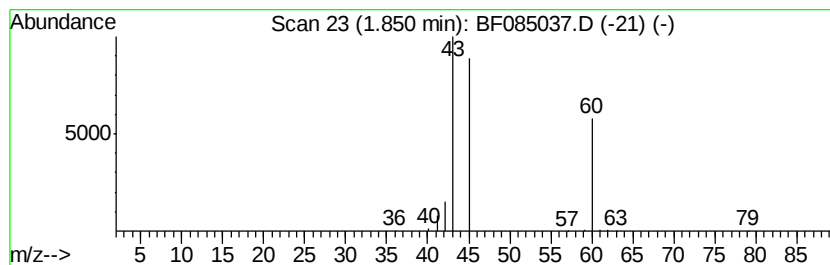
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Peak Number 2 Acetic acid Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.85	220.71 ng	9532450	1,4-Dichlorobenzene-d4	6.60

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Acetic acid		60	C2H4O2	000064-19-7	90
2	Acetic acid, anhydride with form...		88	C3H4O3	002258-42-6	9
3	Ammonium acetate		77	C2H7NO2	000631-61-8	9
4	Hydrazine, 1,2-dimethyl-		60	C2H8N2	000540-73-8	5
5	Thiirane		60	C2H4S	000420-12-2	5



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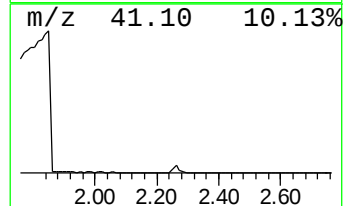
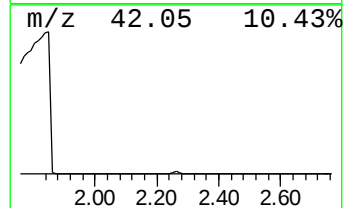
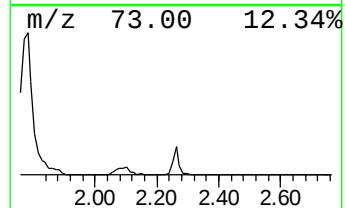
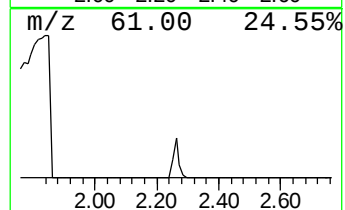
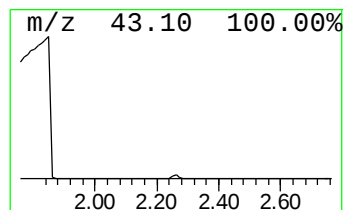
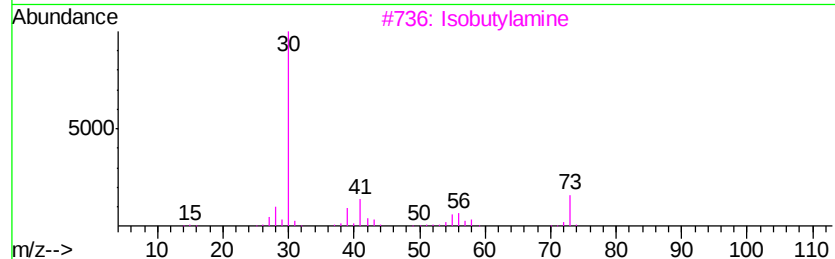
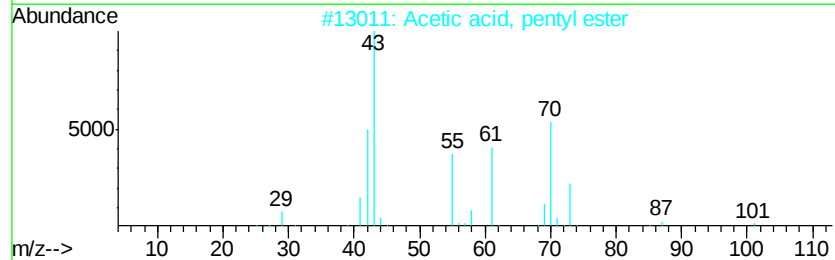
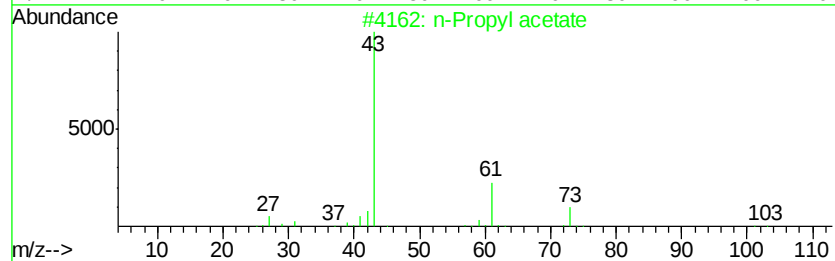
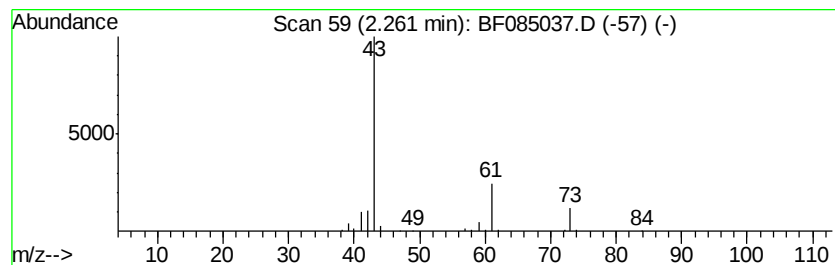
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Peak Number 3 n-Propyl acetate Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.26	3.68 ng	158995	1,4-Dichlorobenzene-d4	6.60

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Propyl acetate	102	C5H10O2	000109-60-4	83
2		Acetic acid, pentyl ester	130	C7H14O2	000628-63-7	25
3		Isobutylamine	73	C4H11N	000078-81-9	5
4		1-Butanamine	73	C4H11N	000109-73-9	5
5		Butane, 2-methyl-	72	C5H12	000078-78-4	4



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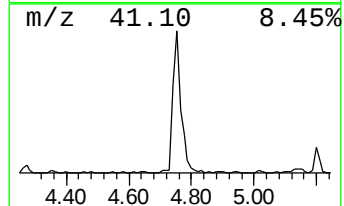
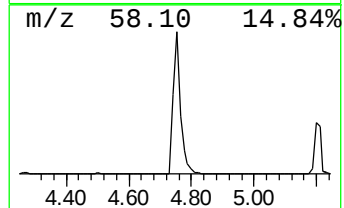
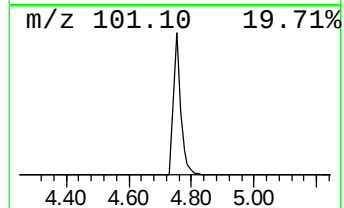
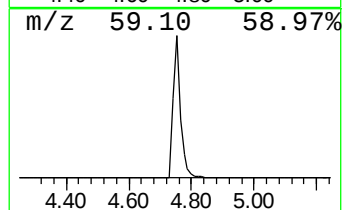
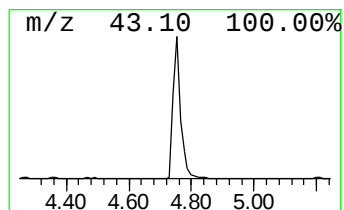
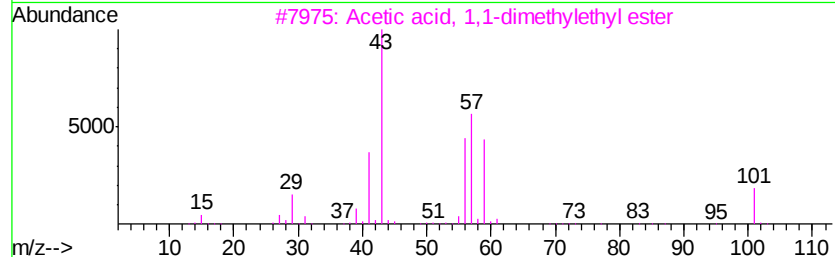
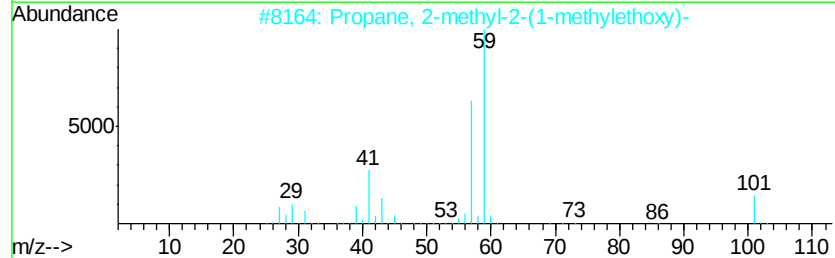
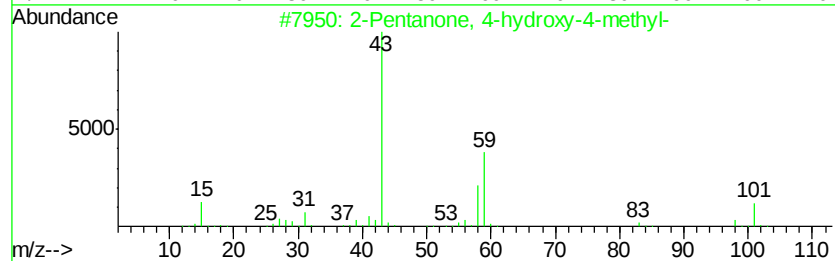
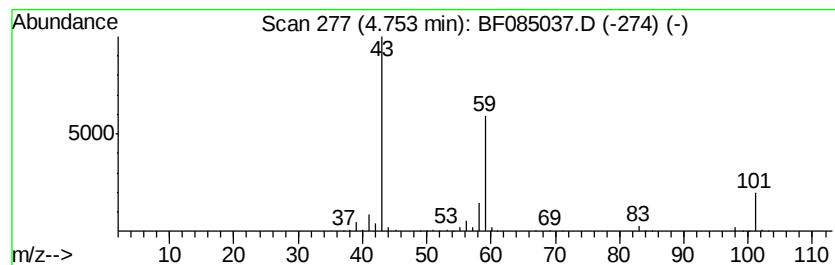
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Peak Number 4 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.75	31.19 ng	1347100	1,4-Dichlorobenzene-d4	6.60

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	72
2		Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	53
3		Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
4		3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	28
5		Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	23



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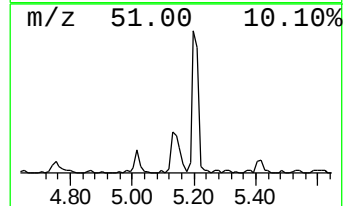
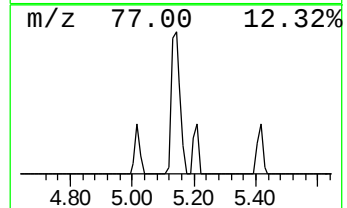
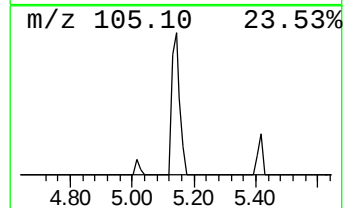
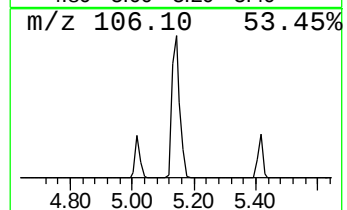
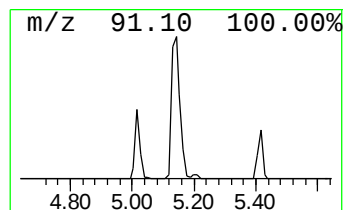
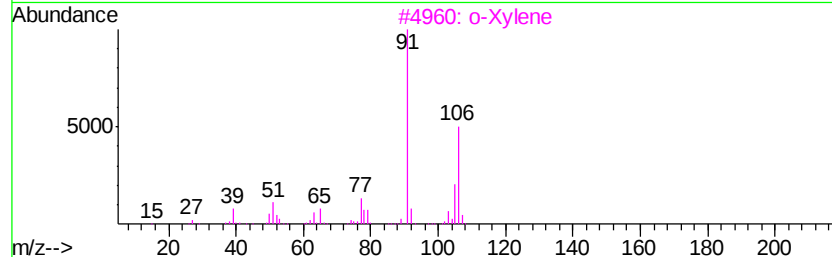
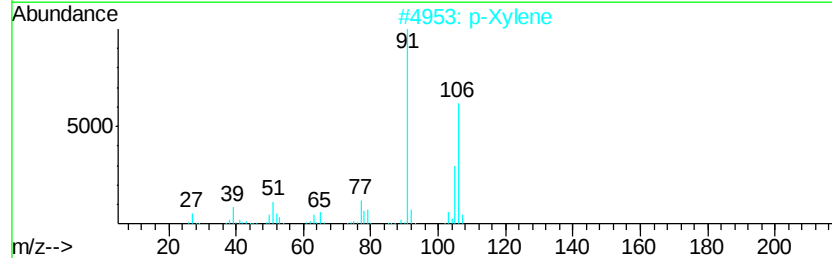
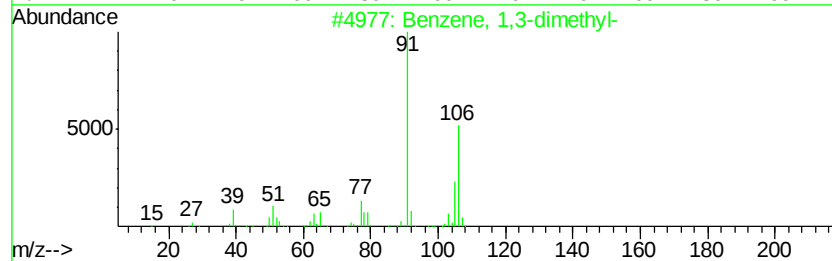
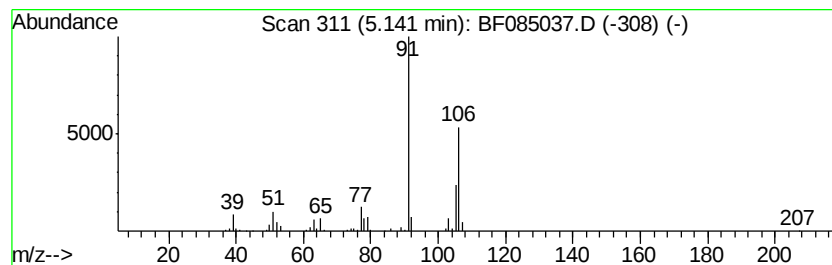
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Peak Number 5 Benzene, 1,3-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.14	7.85 ng	338989	1,4-Dichlorobenzene-d4	6.60

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	97
2		p-Xylene	106	C8H10	000106-42-3	95
3		o-Xylene	106	C8H10	000095-47-6	94
4		1,3-Cyclopentadiene, 5-(1-methyl...	106	C8H10	002175-91-9	87
5		Cyclopentene, 1-ethenyl-3-methyl...	106	C8H10	061142-07-2	86



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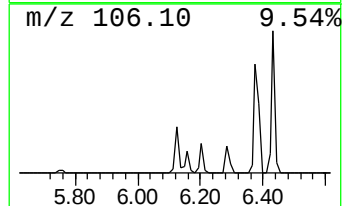
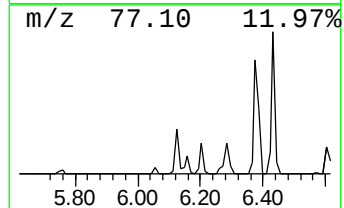
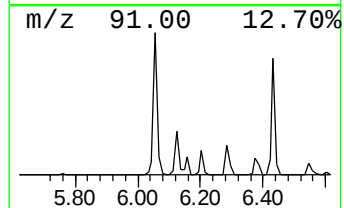
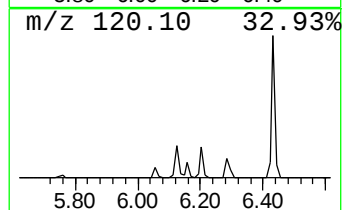
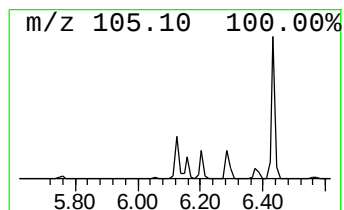
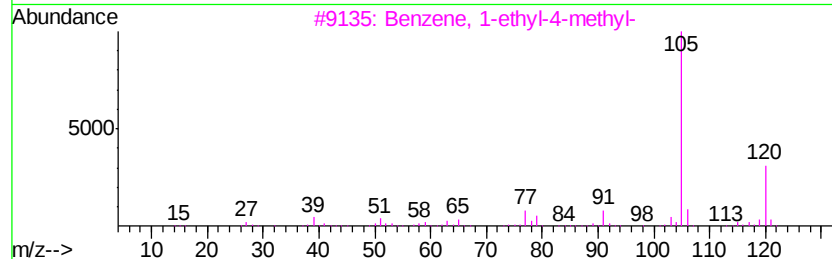
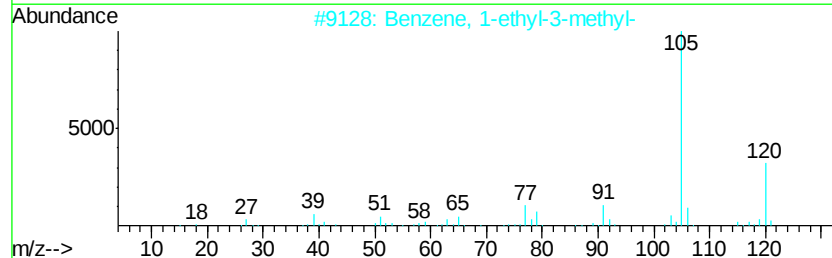
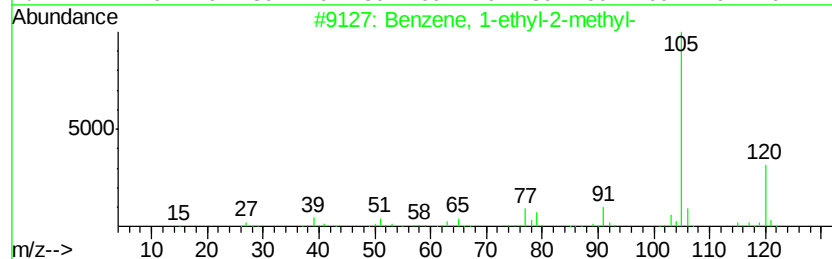
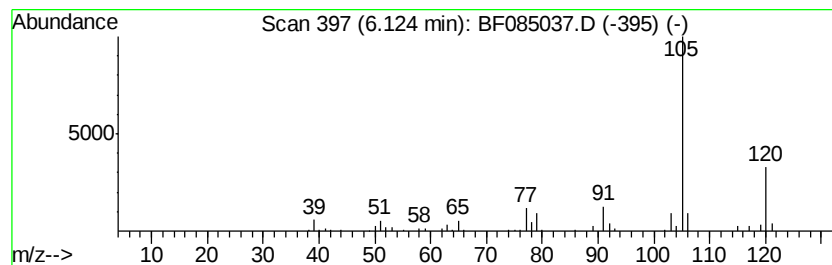
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Peak Number 6 Benzene, 1-ethyl-2-methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.12	3.58 ng	154741	1,4-Dichlorobenzene-d4	6.60

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
2		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	95
3		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
4		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
5		Benzene, 1,3,5-trimethyl-	120	C9H12	000108-67-8	91



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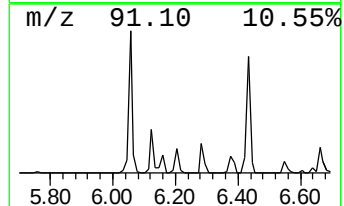
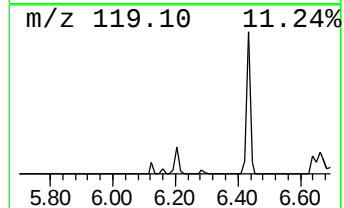
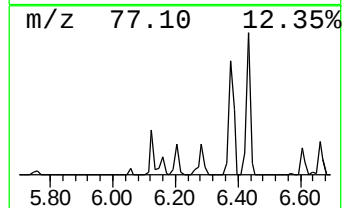
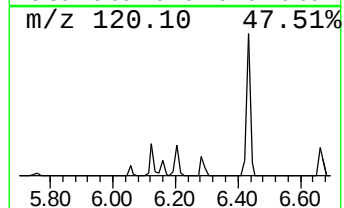
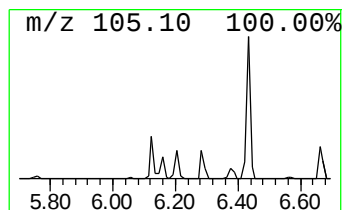
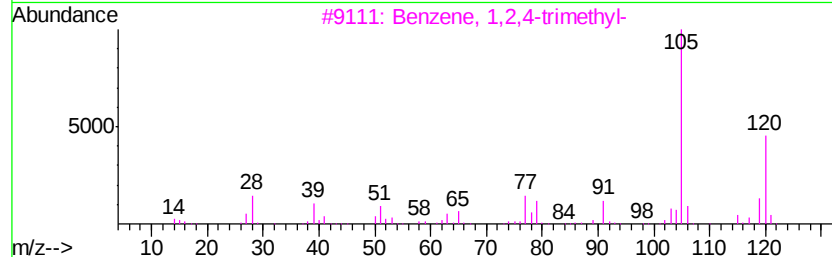
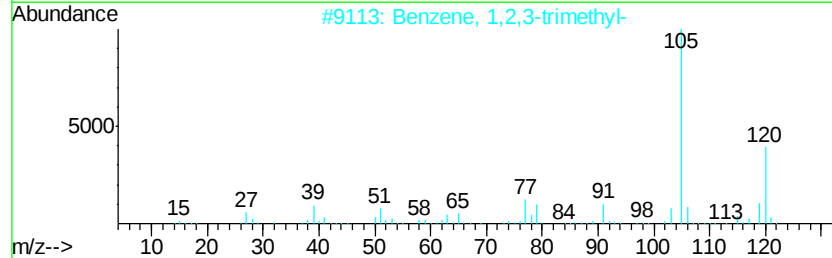
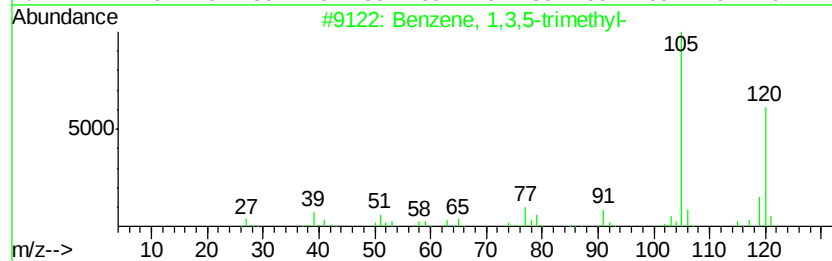
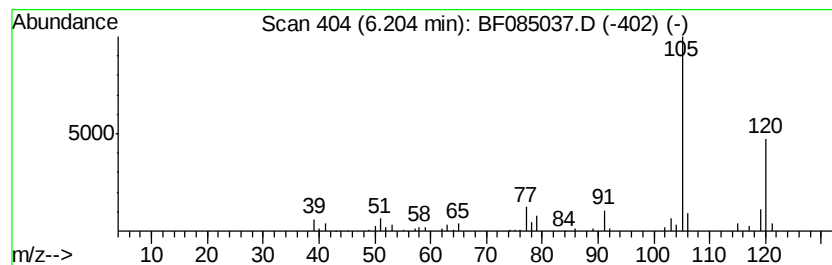
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Benzene, 1,3,5-trimethyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.20	2.77 ng	119453	1,4-Dichlorobenzene-d4	6.60

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,3,5-trimethyl-	120	C9H12	000108-67-8	97
2		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
3		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
4		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	90
5		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	90



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF021216\
Data File : BF085037.D
Acq On : 12 Feb 2016 3:16
Operator : SJ/IZ
Sample : H1410-01
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SUP-B004(16-18)

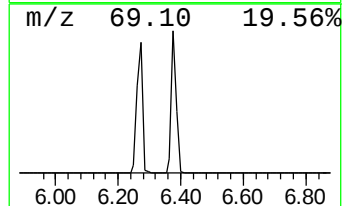
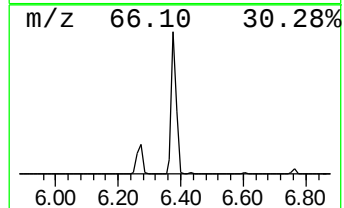
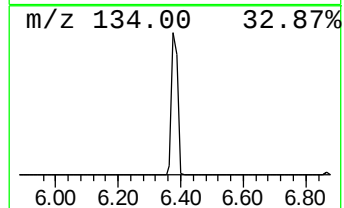
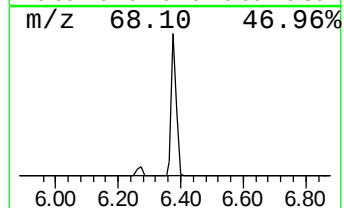
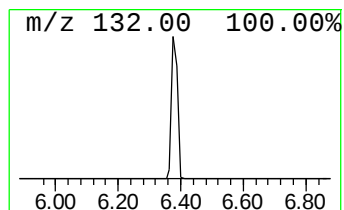
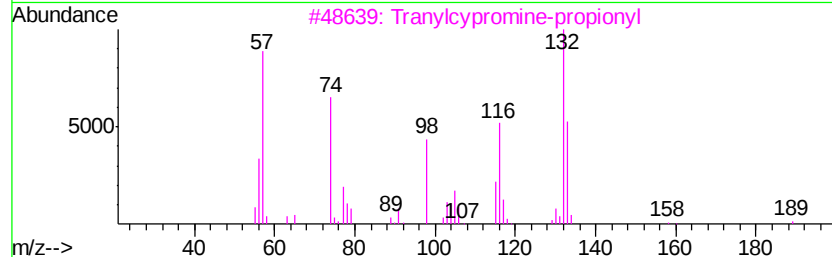
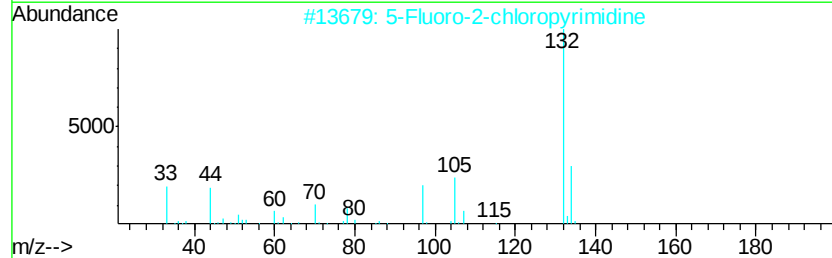
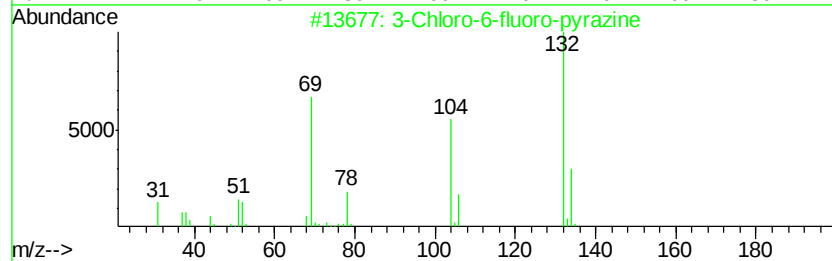
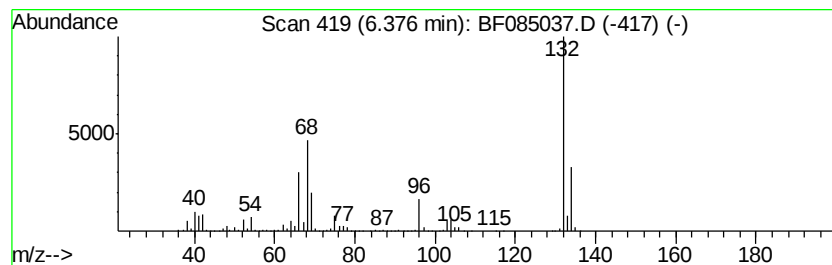
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 unknown6.38 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.38	92.69 ng	4003460	1,4-Dichlorobenzene-d4	6.60

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
3		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
4		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
5		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF021216\
Data File : BF085037.D
Acq On : 12 Feb 2016 3:16
Operator : SJ/IZ
Sample : H1410-01
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
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ClientSampleId :
SUP-B004(16-18)

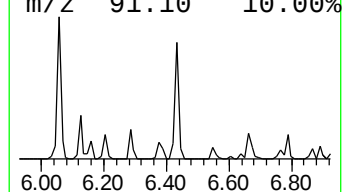
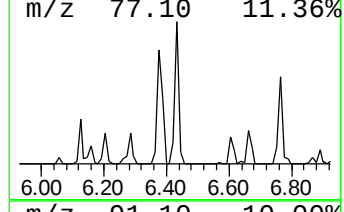
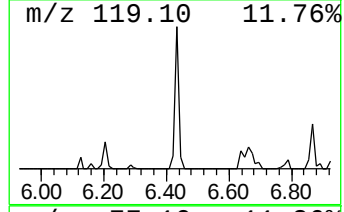
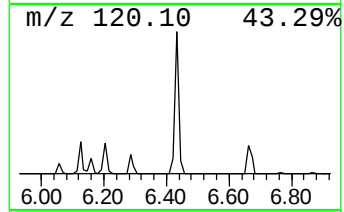
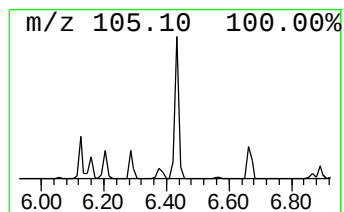
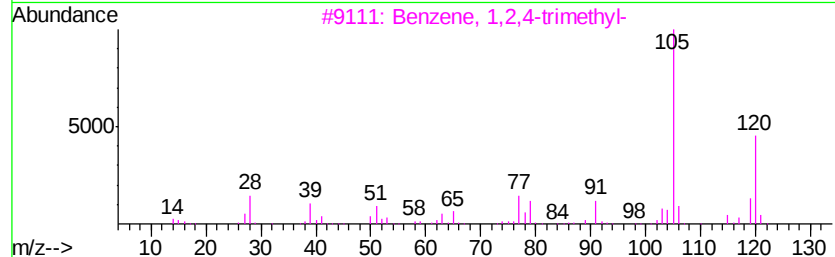
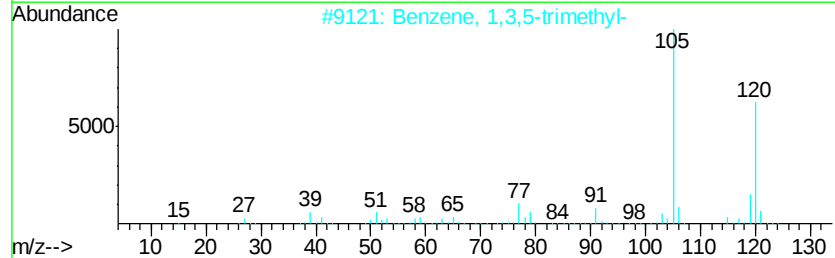
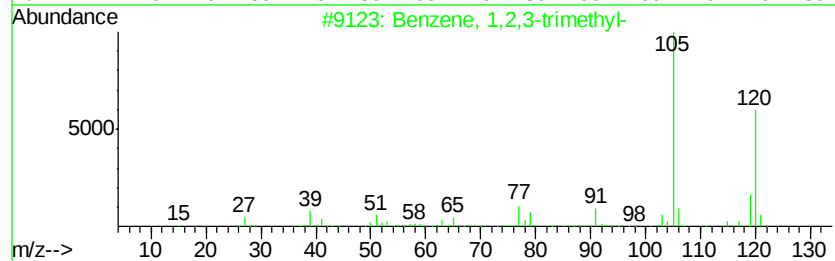
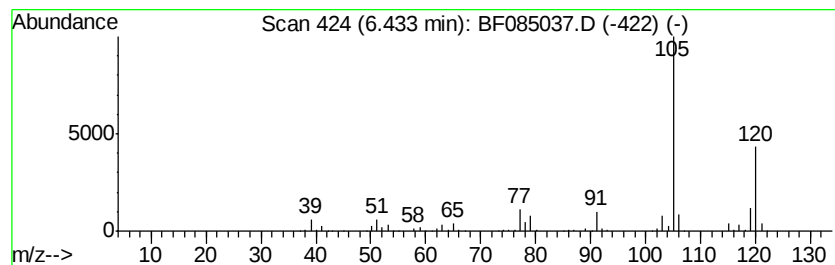
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Benzene, 1,2,3-trimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.43	13.56 ng	585452	1,4-Dichlorobenzene-d4	6.60

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	97
2		Benzene, 1,3,5-trimethyl-	120	C9H12	000108-67-8	94
3		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
4		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	90
5		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	87



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF021216\
Data File : BF085037.D
Acq On : 12 Feb 2016 3:16
Operator : SJ/IZ
Sample : H1410-01
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
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ClientSampleId :
SUP-B004(16-18)

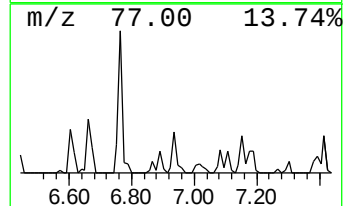
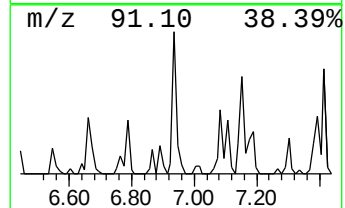
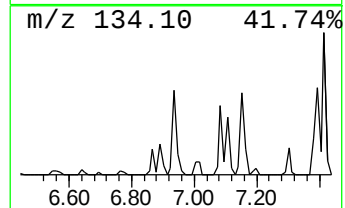
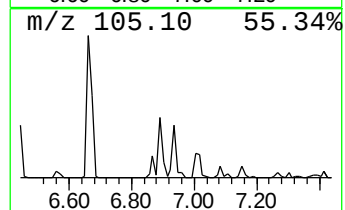
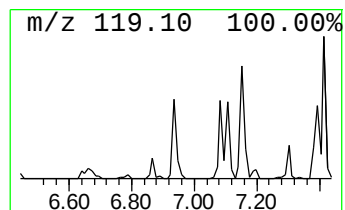
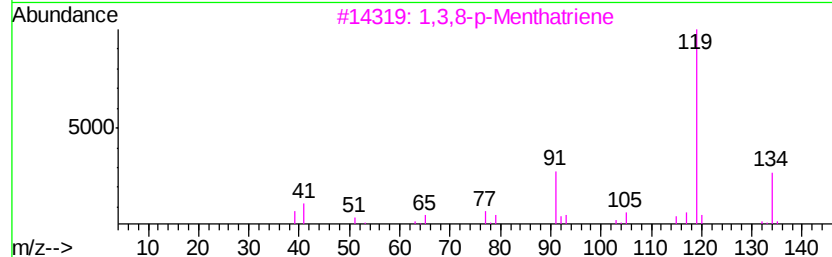
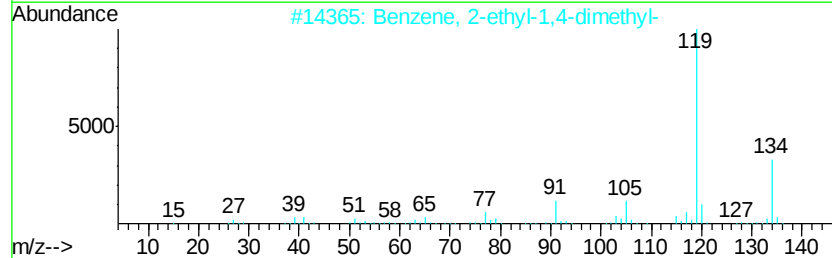
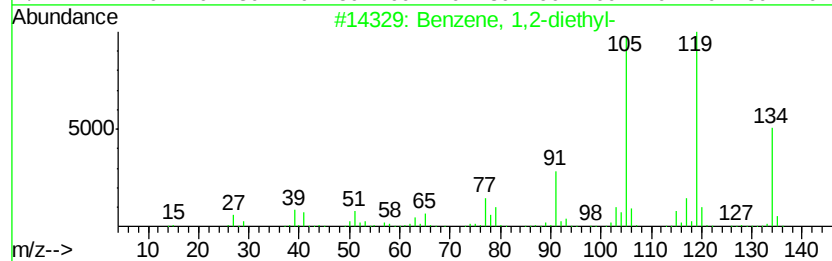
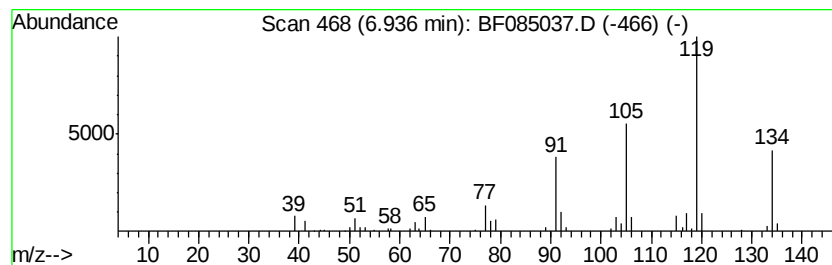
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Benzene, 1,2-diethyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.94	3.20 ng	138293	1,4-Dichlorobenzene-d4	6.60

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	92
2			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	91
3			1,3,8-p-Menthatriene	134	C10H14	021195-59-5	91
4			Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	91
5			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	91



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF021216\
Data File : BF085037.D
Acq On : 12 Feb 2016 3:16
Operator : SJ/IZ
Sample : H1410-01
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SUP-B004(16-18)

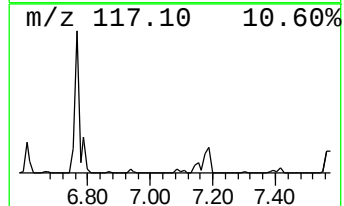
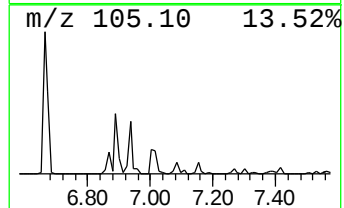
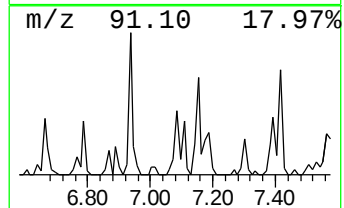
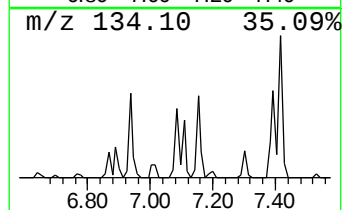
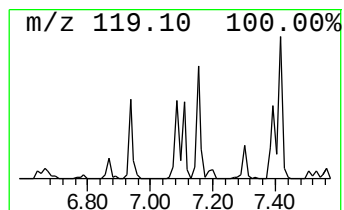
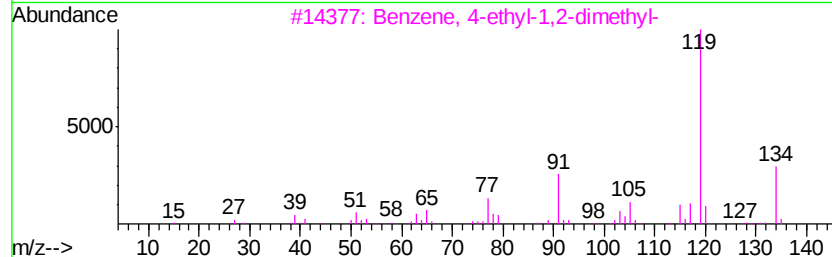
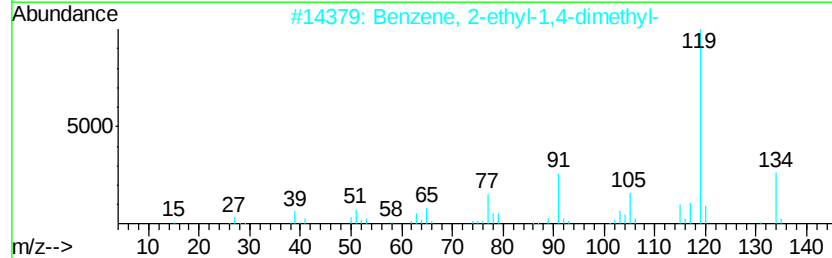
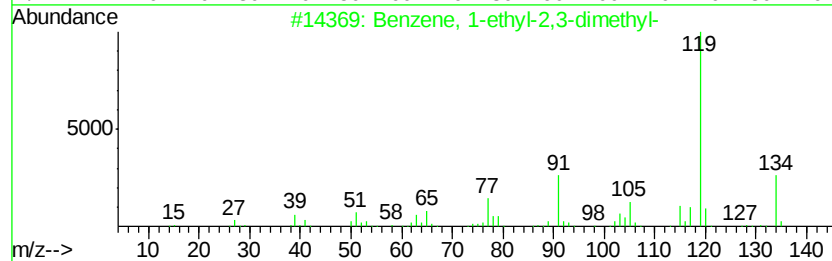
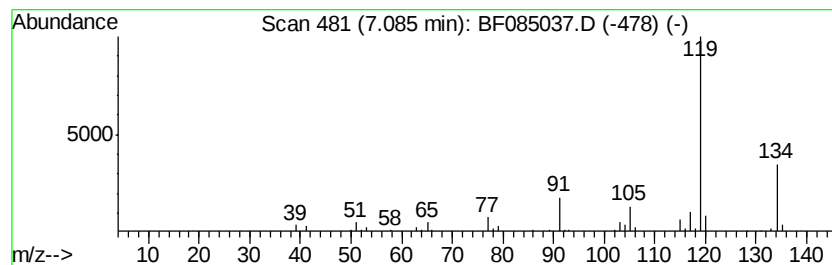
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Benzene, 1-ethyl-2,3-dimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.08	3.74 ng	161727	1,4-Dichlorobenzene-d4	6.60

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	96
2		Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	96
3		Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
4		Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	91
5		Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	91



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF021216\
Data File : BF085037.D
Acq On : 12 Feb 2016 3:16
Operator : SJ/IZ
Sample : H1410-01
Misc :
ALS Vial : 27 Sample Multiplier: 1

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ClientSampleId :
SUP-B004(16-18)

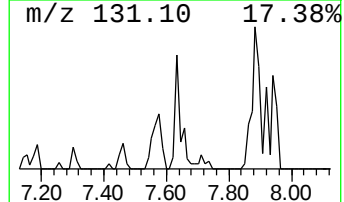
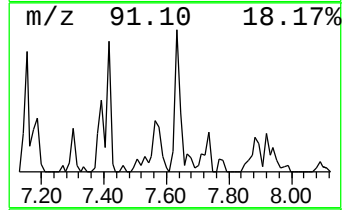
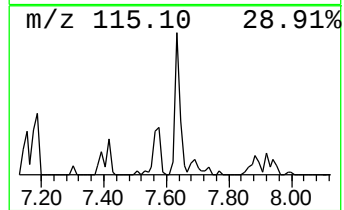
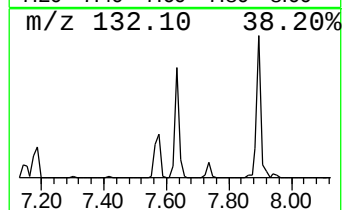
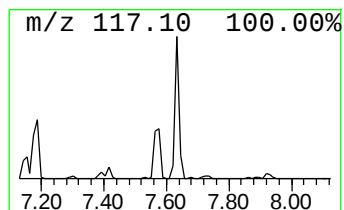
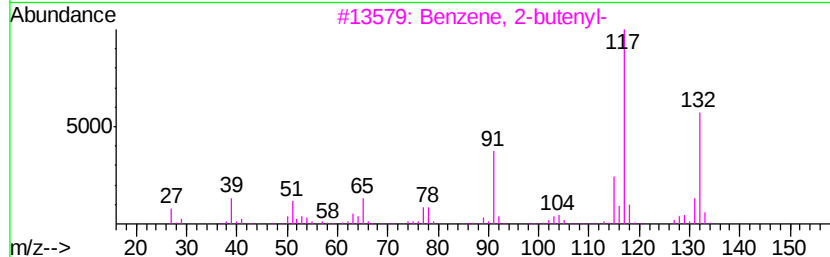
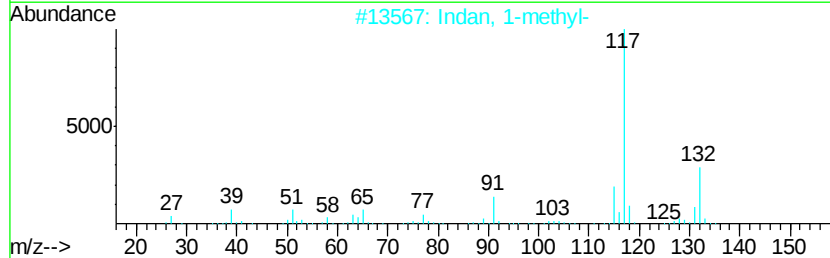
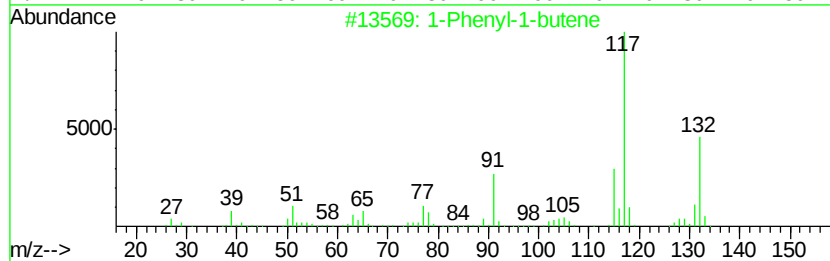
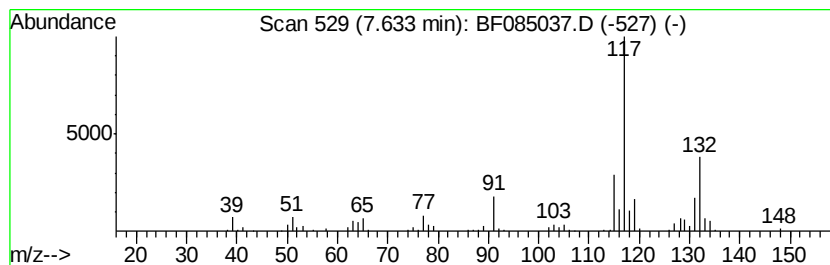
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 1-Phenyl-1-butene Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.63	3.38 ng	287016	Naphthalene-d8	7.90

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Phenyl-1-butene	132	C10H12	000824-90-8	93
2		Indan, 1-methyl-	132	C10H12	000767-58-8	90
3		Benzene, 2-butenyl-	132	C10H12	001560-06-1	90
4		Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	87
5		1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	87



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF021216\
Data File : BF085037.D
Acq On : 12 Feb 2016 3:16
Operator : SJ/IZ
Sample : H1410-01
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SUP-B004(16-18)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF020116.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
Acetic acid	1.85	220.7	ng	9532450	1	6.60	863813	20.0
n-Propyl acetate	2.26	3.7	ng	158995	1	6.60	863813	20.0
2-Pentanone, 4-hy...	4.75	31.2	ng	1347100	1	6.60	863813	20.0
Benzene, 1,3-dime...	5.14	7.8	ng	338989	1	6.60	863813	20.0
Benzene, 1-ethyl-...	6.12	3.6	ng	154741	1	6.60	863813	20.0
Benzene, 1,3,5-tr...	6.20	2.8	ng	119453	1	6.60	863813	20.0
unknown6.38	6.38	92.7	ng	4003460	1	6.60	863813	20.0
Benzene, 1,2,3-tr...	6.43	13.6	ng	585452	1	6.60	863813	20.0
Benzene, 1,2-diet...	6.94	3.2	ng	138293	1	6.60	863813	20.0
Benzene, 1-ethyl-...	7.08	3.7	ng	161727	1	6.60	863813	20.0
1-Phenyl-1-butene	7.63	3.4	ng	287016	2	7.90	1695890	20.0