

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG080818\
 Data File : BG036222.D
 Acq On : 9 Aug 2018 7:51
 Operator : JU/SJ
 Sample : J4328-01
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 MW-22

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:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG071218.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	4.639	225	230	235	rBV2	23575	40709	1.47%	0.333%
2	5.126	308	313	319	rBV2	22613	41809	1.51%	0.342%
3	5.596	385	393	406	rBV	193880	365310	13.22%	2.986%
4	7.182	657	663	674	rBV2	110509	240571	8.71%	1.967%
5	7.546	716	725	740	rVB	417998	755973	27.36%	6.180%
6	8.011	796	804	810	rBV	103062	183216	6.63%	1.498%
7	8.334	851	859	870	rBV	392087	724375	26.22%	5.922%
8	9.174	995	1002	1020	rVB	315740	653850	23.67%	5.345%
9	9.368	1028	1035	1041	rVB8	14253	30154	1.09%	0.247%
10	9.514	1053	1060	1068	rVB	166343	283719	10.27%	2.319%
11	9.785	1102	1106	1115	rVB4	13628	28018	1.01%	0.229%
12	10.208	1173	1178	1183	rBV4	18020	33830	1.22%	0.277%
13	10.813	1274	1281	1294	rVB	118115	253618	9.18%	2.073%
14	10.965	1302	1307	1312	rVB2	19540	31806	1.15%	0.260%
15	11.923	1464	1470	1477	rVB8	13764	29276	1.06%	0.239%
16	13.256	1690	1697	1709	rBV	917432	1500906	54.32%	12.270%
17	13.814	1785	1792	1800	rBV2	22914	49899	1.81%	0.408%
18	13.950	1809	1815	1822	rBV3	32966	64488	2.33%	0.527%
19	14.190	1850	1856	1860	rBV4	28602	52989	1.92%	0.433%
20	14.631	1925	1931	1937	rBV2	222991	372988	13.50%	3.049%
21	14.696	1937	1942	1950	rVB	139465	239791	8.68%	1.960%
22	14.842	1961	1967	1974	rBV	13754	31634	1.14%	0.259%
23	15.030	1993	1999	2007	rVB	76144	131224	4.75%	1.073%
24	15.677	2105	2109	2115	rVB2	36969	58792	2.13%	0.481%
25	16.123	2179	2185	2194	rBV	838515	1370235	49.59%	11.201%
26	16.646	2269	2274	2279	rBV	60566	102335	3.70%	0.837%
27	17.374	2393	2398	2405	rBV	323425	498620	18.05%	4.076%
28	18.485	2582	2587	2592	rBV3	44935	79545	2.88%	0.650%
29	19.988	2837	2843	2849	rBV	2162816	2762872	100.00%	22.586%
30	21.639	3117	3124	3134	rBV2	360634	633181	22.92%	5.176%
31	24.823	3655	3666	3678	rBV2	187550	586979	21.25%	4.798%

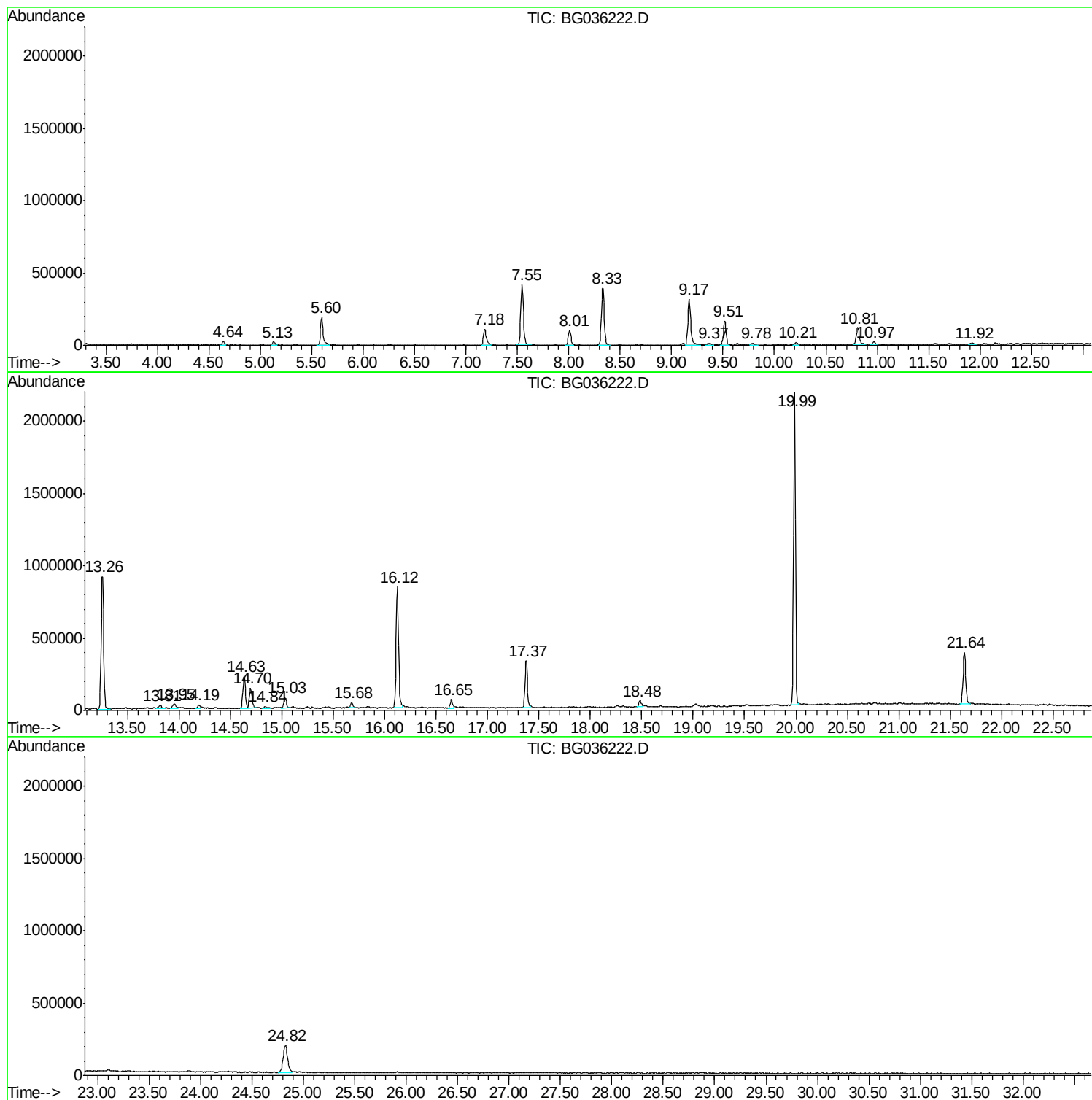
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG071218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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Instrument :
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ClientSampled :
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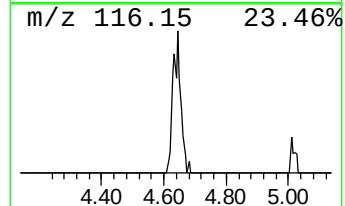
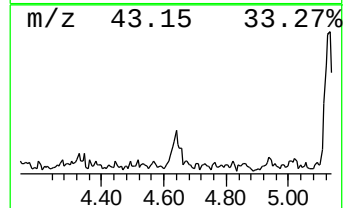
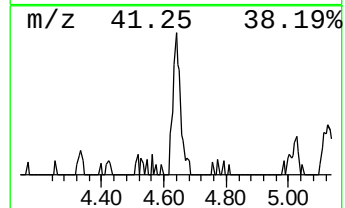
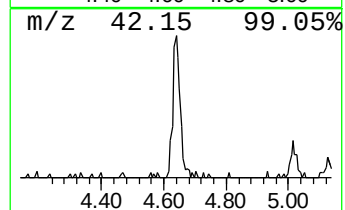
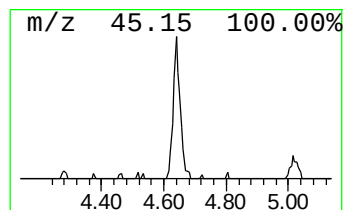
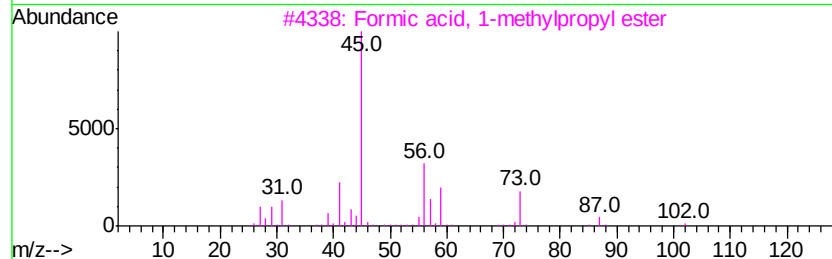
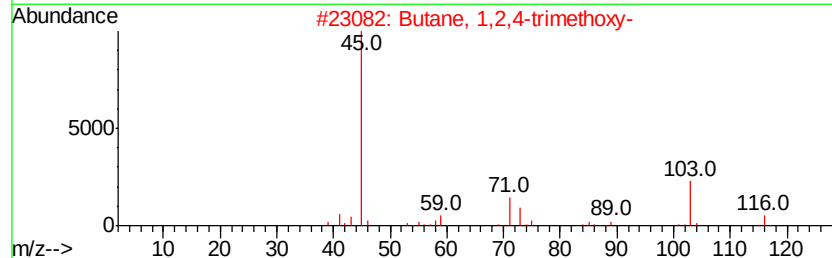
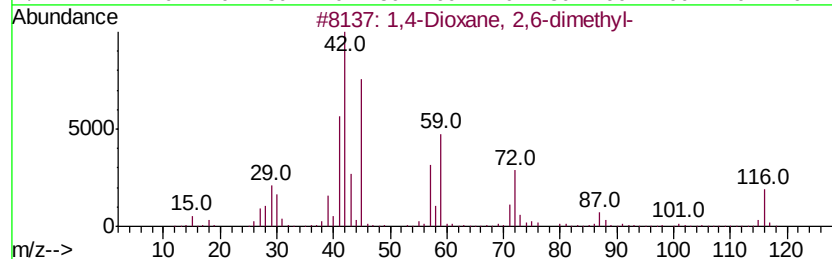
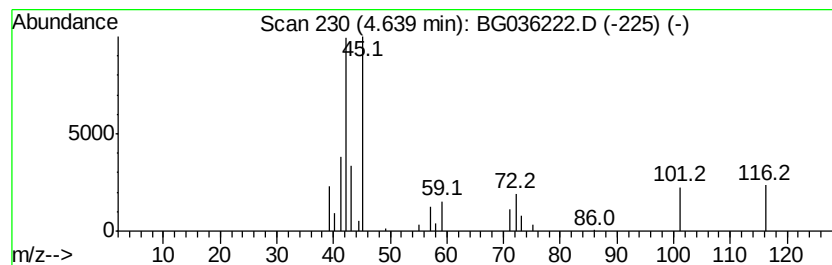
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG071218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 unknown4.64 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.64	4.44 ng	40709	1,4-Dichlorobenzene-d4	8.01

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,4-Dioxane, 2,6-dimethyl-	116	C6H12O2	010138-17-7	32
2		Butane, 1,2,4-trimethoxy-	148	C7H16O3	020637-48-3	12
3		Formic acid, 1-methylpropyl ester	102	C5H10O2	000589-40-2	12
4		2,3-Butanedione, dioxime	116	C4H8N2O2	000095-45-4	9
5		2-Methoxyisopropyl cyanoacetate	157	C7H11NO3	032804-79-8	9



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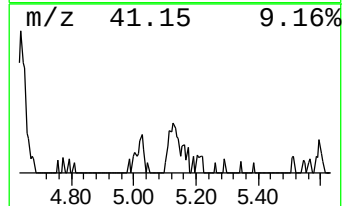
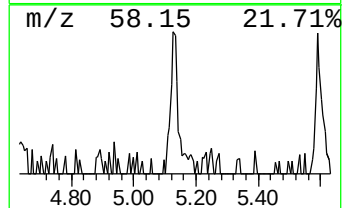
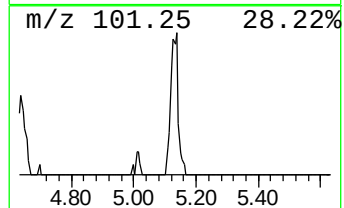
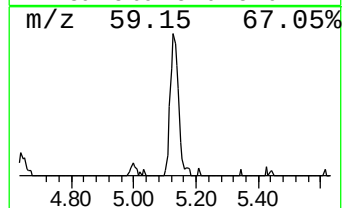
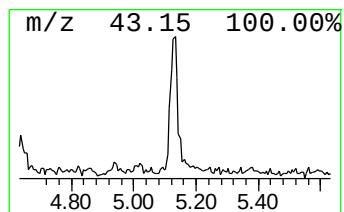
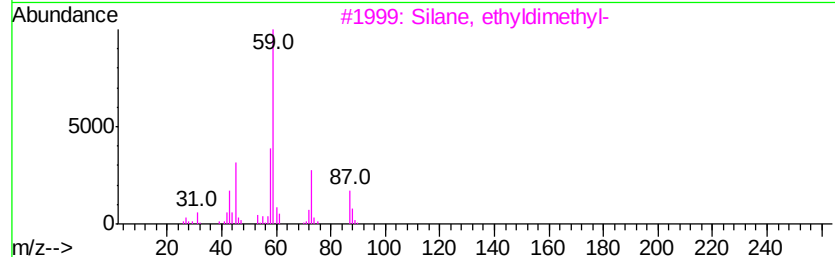
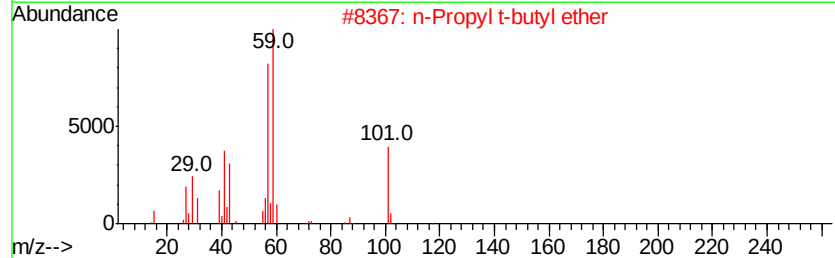
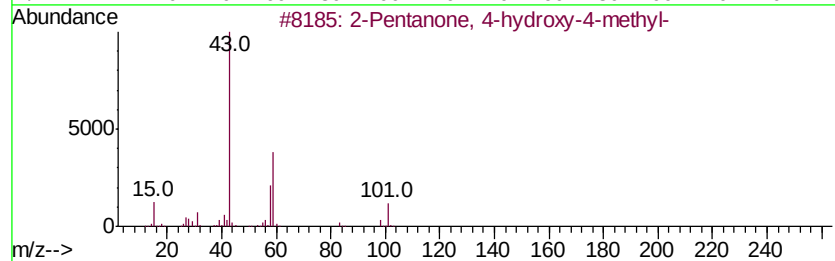
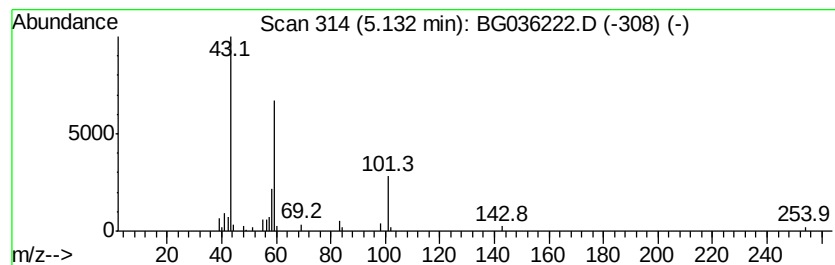
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TIC Library : C:\Database\NIST11.L
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Peak Number 2 unknown5.13 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.13	4.56 ng	41809	1,4-Dichlorobenzene-d4	8.01

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	45
2	n-Propyl t-butyl ether	116	C7H16O	1000334-81-9	42
3	Silane, ethyldimethyl-	88	C4H12Si	000758-21-4	38
4	2-Nonanone, 9-hydroxy-	158	C9H18O2	025368-56-3	28
5	4-Methoxy-1-pentene	100	C6H12O	098386-09-5	25



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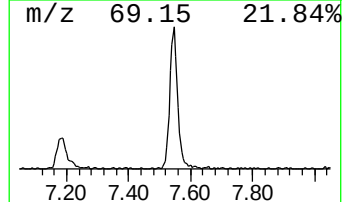
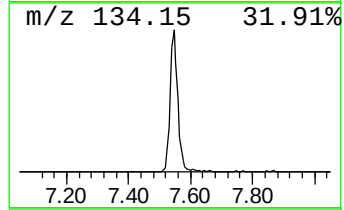
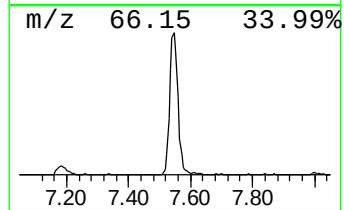
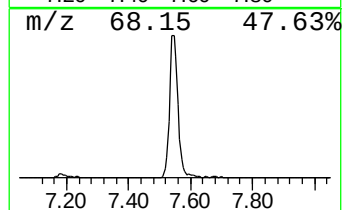
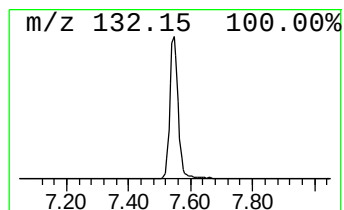
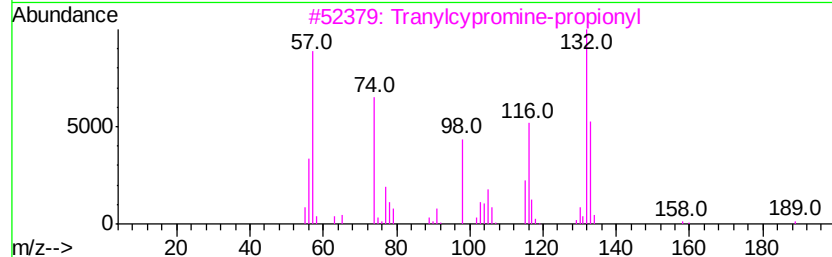
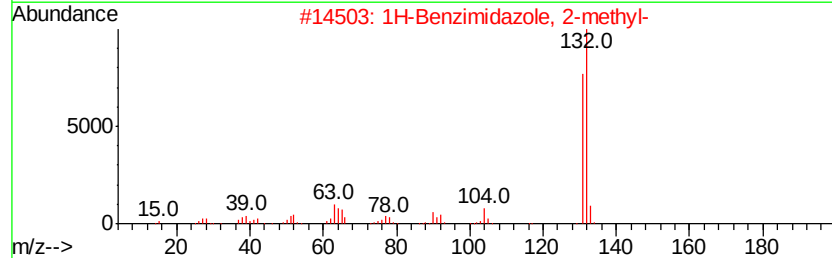
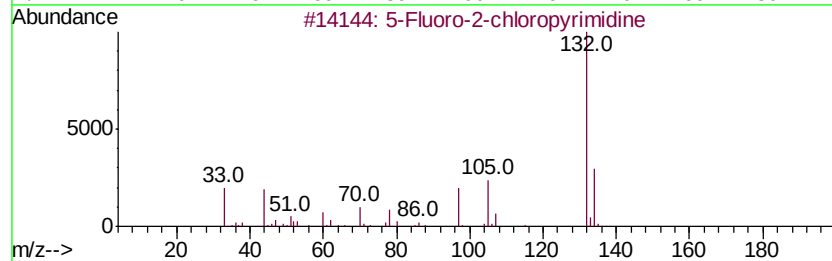
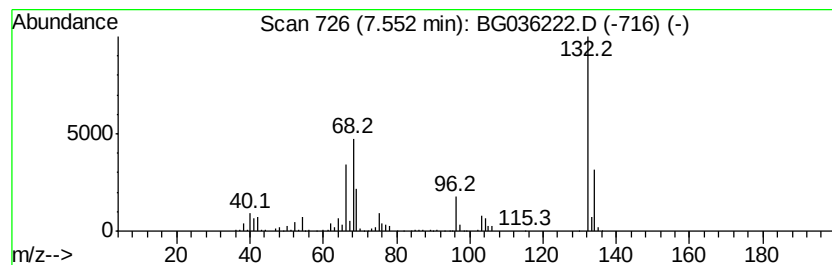
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Peak Number 3 unknown7.55 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.55	82.52 ng	755973	1,4-Dichlorobenzene-d4	8.01

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	27
2		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
3		Tranvlcypromine-propionyl	189	C12H15NO	1000123-86-3	22
4		(5-Methyl-2-pyridyl)acetonitrile	132	C8H8N2	1000241-93-9	22
5		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16



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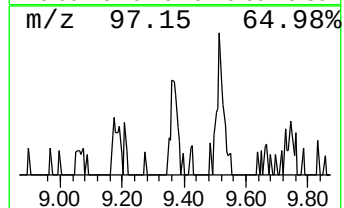
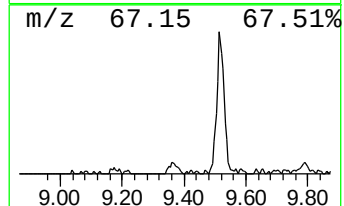
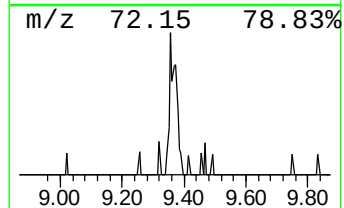
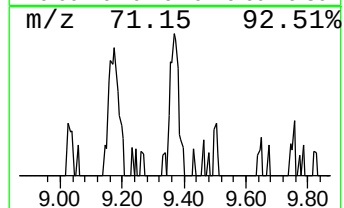
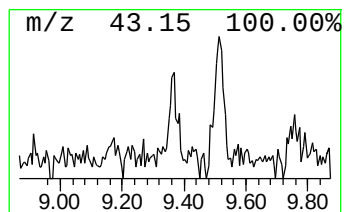
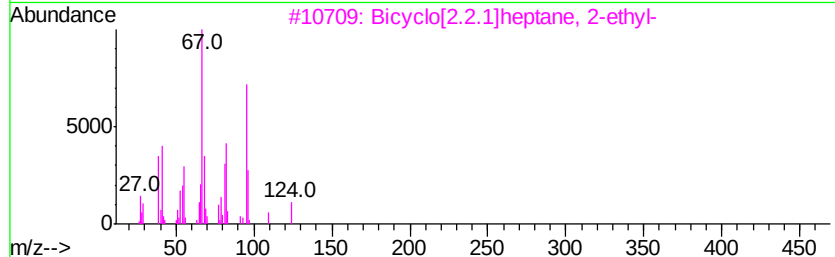
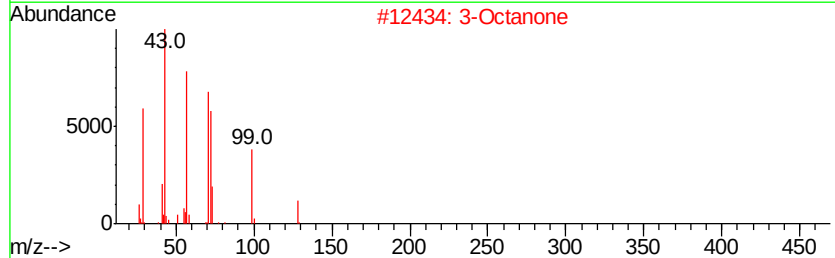
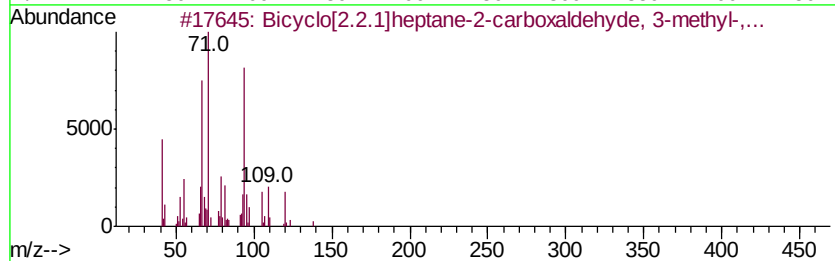
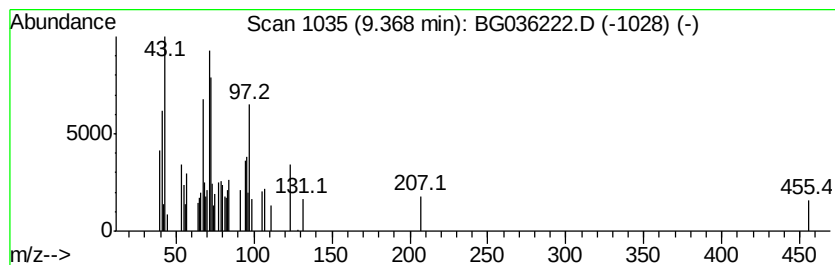
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Peak Number 5 unknown9.37 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.37	3.29 ng	30154	1,4-Dichlorobenzene-d4	8.01

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Bicyclo[2.2.1]heptane-2-carboxal...	138	C9H14O	015780-37-7	32
2		3-Octanone	128	C8H16O	000106-68-3	27
3		Bicyclo[2.2.1]heptane, 2-ethyl-	124	C9H16	002146-41-0	14
4		4,7-Methano-1H-indene, octahydro-	136	C10H16	006004-38-2	14
5		1,2-Cyclohexanedione dioxime	142	C6H10N2O2	000492-99-9	14



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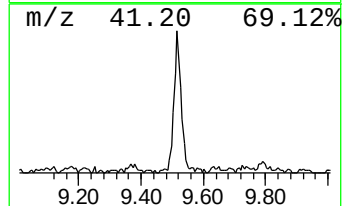
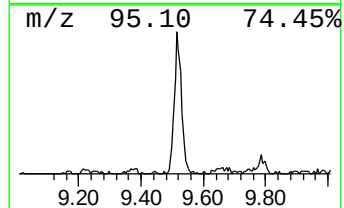
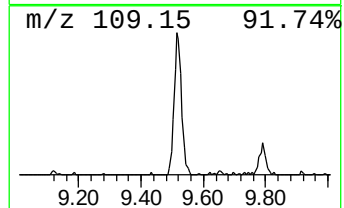
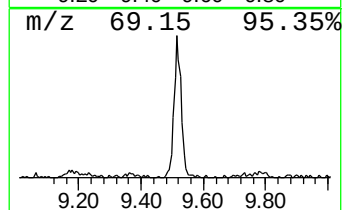
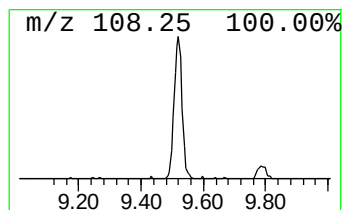
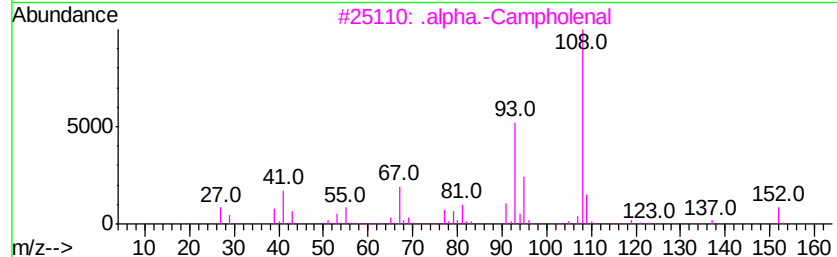
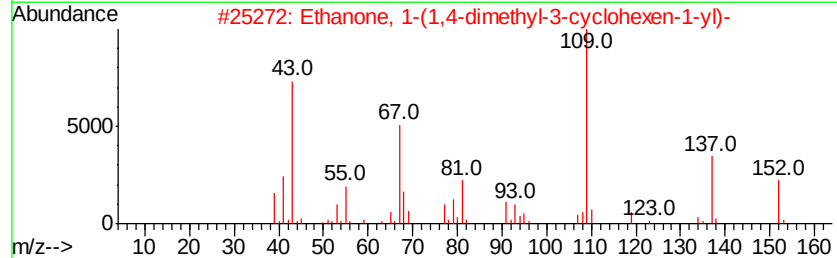
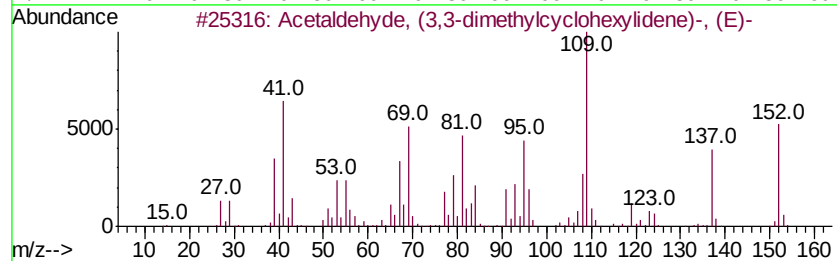
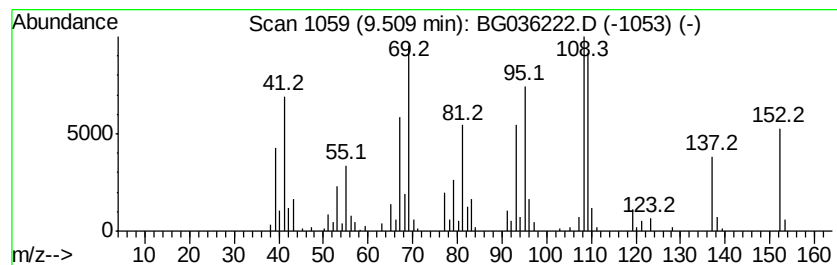
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Peak Number 6 unknown9.51 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.51	22.37 ng	283719	Naphthalene-d8	10.81

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Acetaldehyde, (3,3-dimethylcyclo...	152	C10H16O	026532-25-2	46
2		Ethanone, 1-(1,4-dimethyl-3-cycl...	152	C10H16O	043219-68-7	42
3		.alpha.-Campholenal	152	C10H16O	004501-58-0	27
4		4-Ethyl-2-hexynal	124	C8H12O	071932-97-3	22
5		2,6-Heptadienal, 2,4-dimethyl-	138	C9H14O	085136-08-9	20



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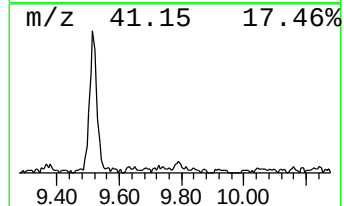
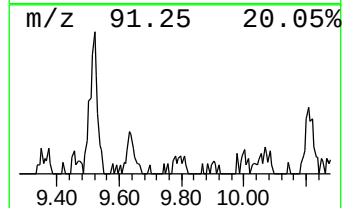
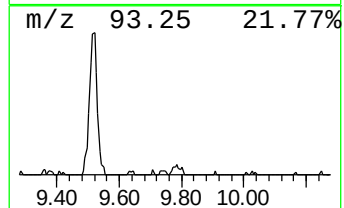
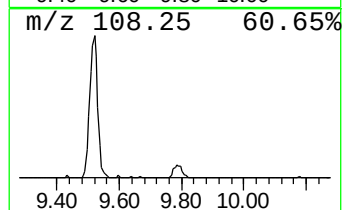
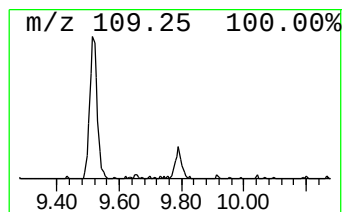
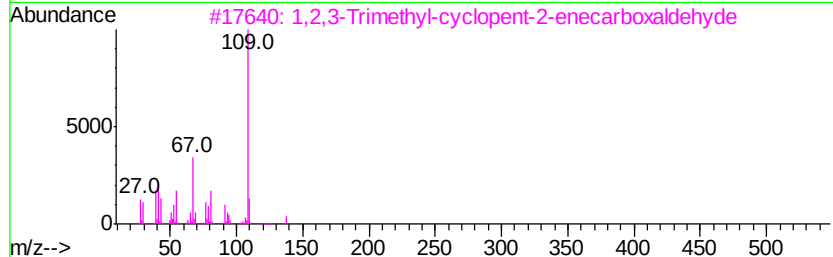
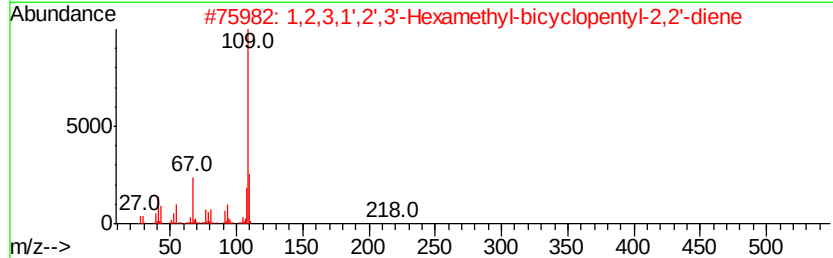
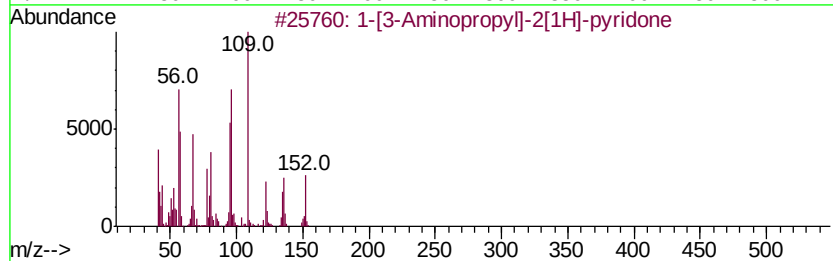
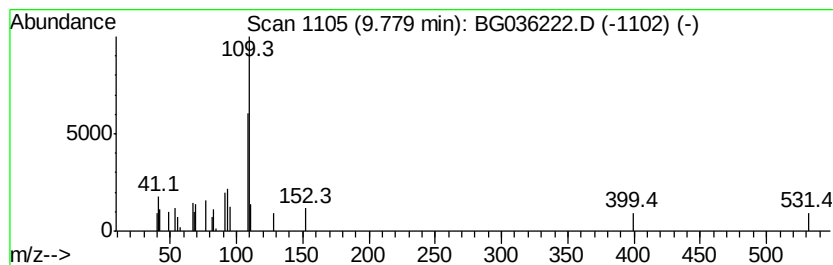
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 unknown9.78 Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.78	2.21 ng	28018	Naphthalene-d8	10.81

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-[3-Aminopropyl]-2[1H]-pyridone	152	C8H12N2O	102675-58-1	47
2		1,2,3,1',2',3'-Hexamethyl-bicycl...	218	C16H26	051999-35-0	43
3		1,2,3-Trimethyl-cyclopent-2-enec...	138	C9H14O	1000190-18-1	38
4		.alpha.-Campholenal	152	C10H16O	004501-58-0	38
5		Phenol, 4-amino-	109	C6H7NO	000123-30-8	35



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG080818\
Data File : BG036222.D
Acq On : 9 Aug 2018 7:51
Operator : JU/SJ
Sample : J4328-01
Misc :
ALS Vial : 26 Sample Multiplier: 1

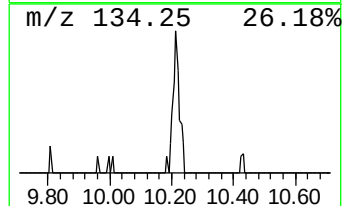
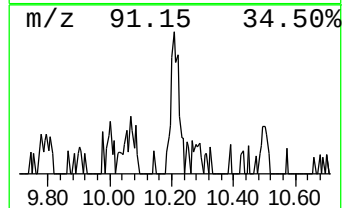
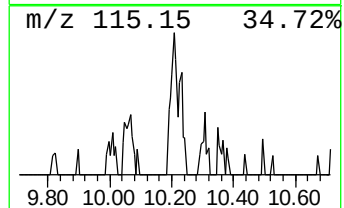
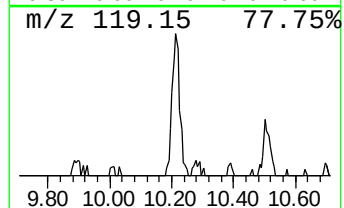
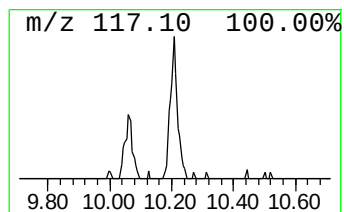
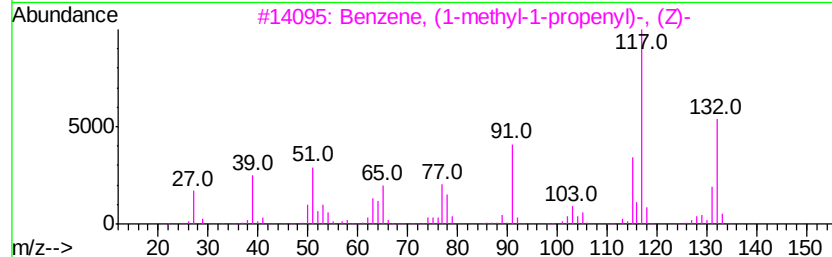
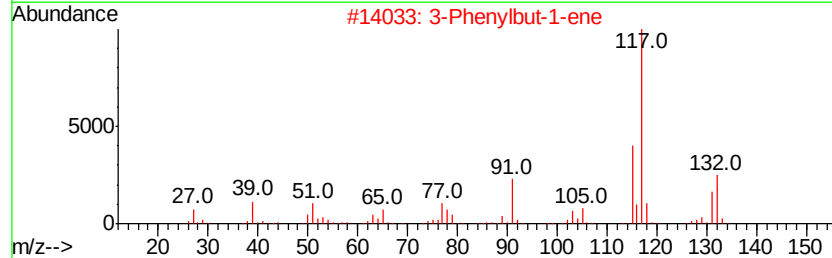
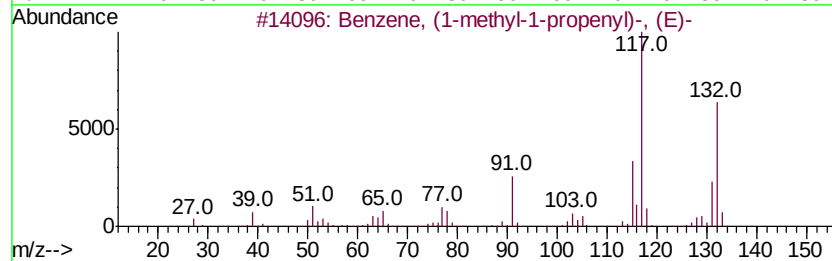
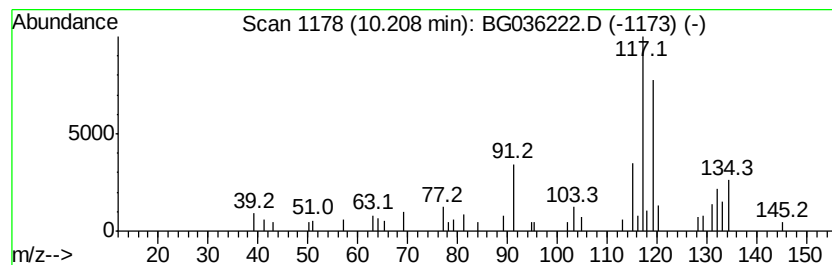
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Benzene, (1-methyl-1-propen... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.21	2.67 ng	33830	Naphthalene-d8	10.81
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, (1-methyl-1-propenyl)-,...	132 C10H12	000768-00-3 55
2		3-Phenylbut-1-ene	132 C10H12	000934-10-1 49
3		Benzene, (1-methyl-1-propenyl)-,...	132 C10H12	000767-99-7 46
4		Benzene, 1-methyl-4-(2-propenyl)-	132 C10H12	003333-13-9 46
5		Benzene, 1-ethenyl-3-ethyl-	132 C10H12	007525-62-4 45



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG080818\
Data File : BG036222.D
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Sample : J4328-01
Misc :
ALS Vial : 26 Sample Multiplier: 1

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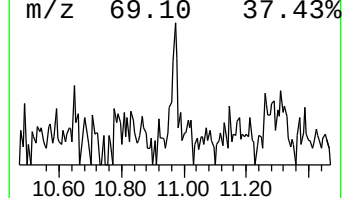
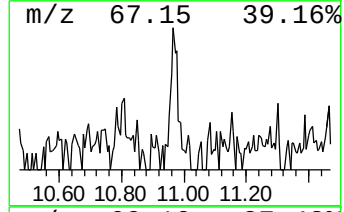
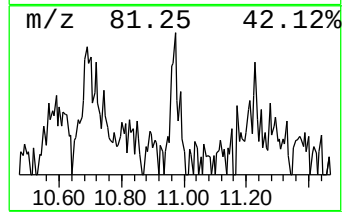
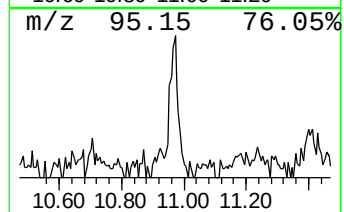
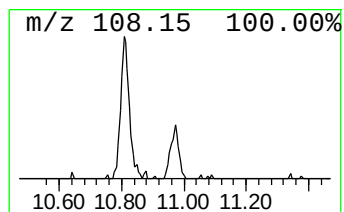
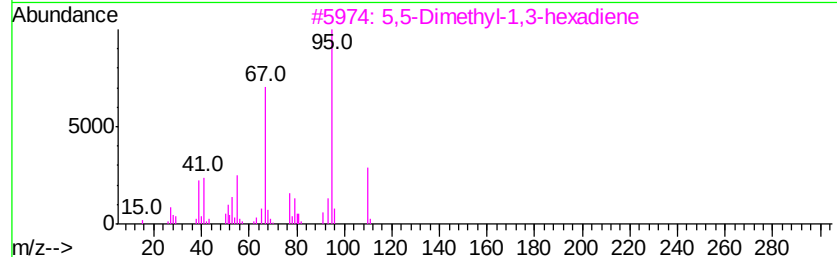
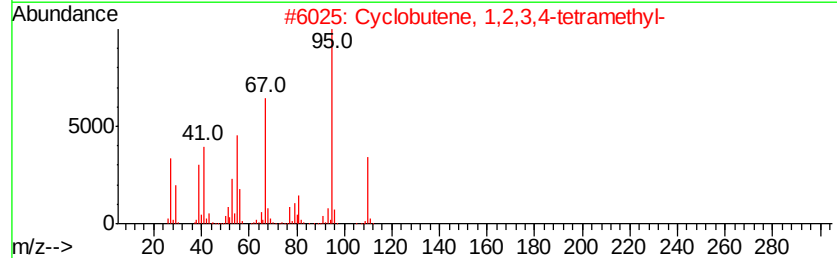
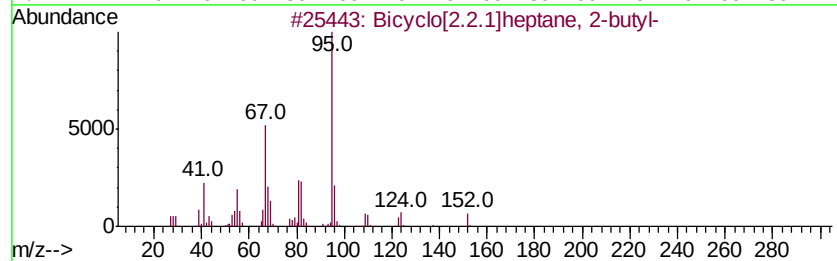
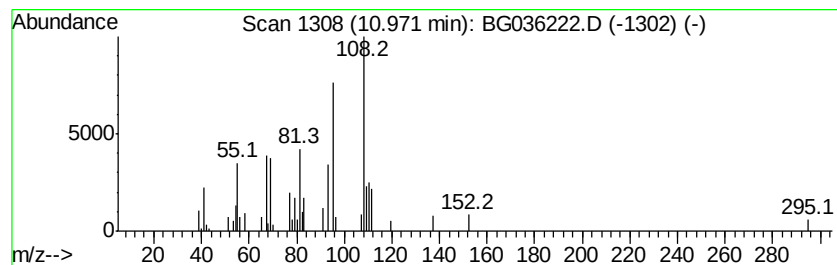
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 unknown10.97 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.97	2.51 ng	31806	Naphthalene-d8	10.81

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Bicyclo[2.2.1]heptane, 2-butyl-	152	C11H20	061177-16-0	35
2		Cyclobutene, 1,2,3,4-tetramethyl-	110	C8H14	090346-45-5	30
3		5,5-Dimethyl-1,3-hexadiene	110	C8H14	001515-79-3	30
4		4-Methyl-1,3-heptadiene	110	C8H14	017603-57-5	30
5		1,3-Hexadiene, 2,5-dimethyl-	110	C8H14	029253-64-3	30



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG080818\
Data File : BG036222.D
Acq On : 9 Aug 2018 7:51
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Sample : J4328-01
Misc :
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ClientSampleId :
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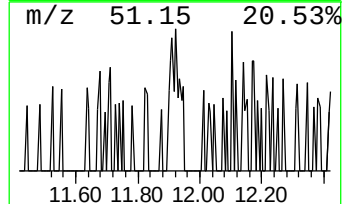
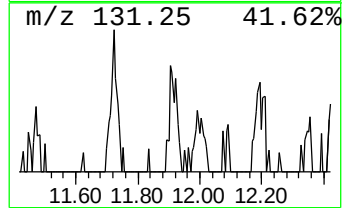
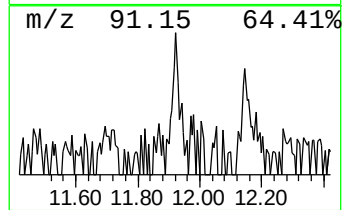
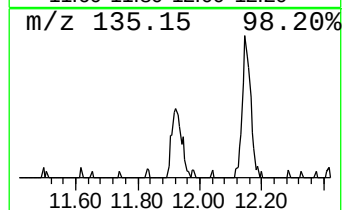
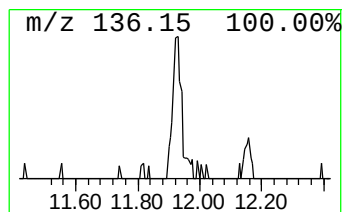
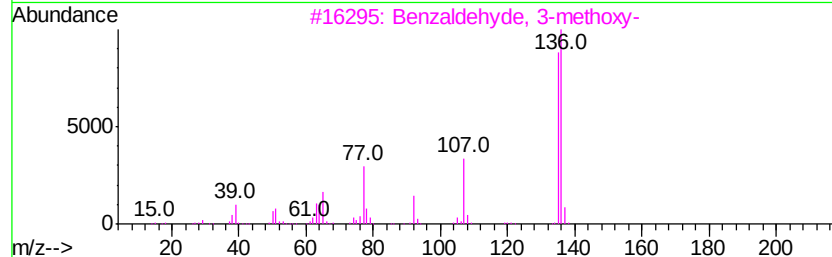
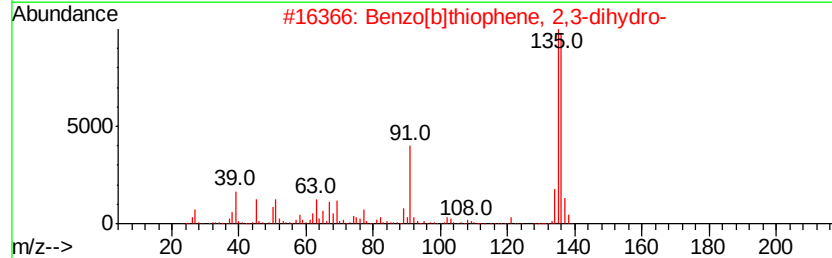
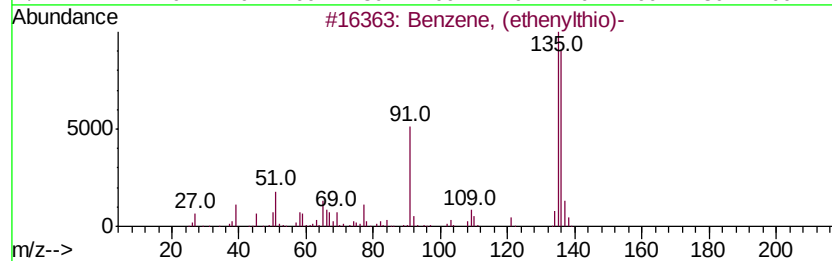
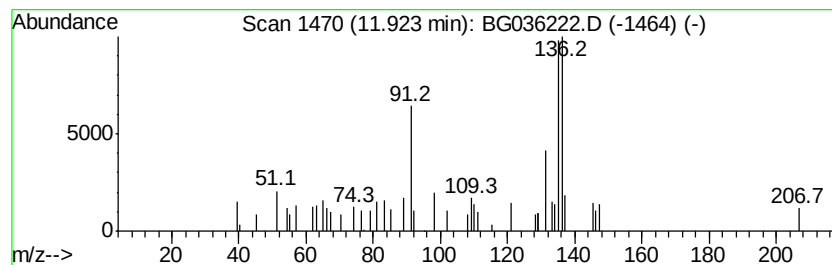
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Benzene, (ethenylthio)- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.92	2.31 ng	29276	Napthalene-d8	10.81

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, (ethenylthio)-	136	C8H8S	001822-73-7	55
2		Benzo[b]thiophene, 2,3-dihydro-	136	C8H8S	004565-32-6	43
3		Benzaldehyde, 3-methoxy-	136	C8H8O2	000591-31-1	43
4		Benzaldehyde, 4-methoxy-	136	C8H8O2	000123-11-5	43
5		3-Furonitrile, 2-amino-4,5-dimet...	136	C7H8N2O	005117-88-4	43



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG080818\
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Sample : J4328-01
Misc :
ALS Vial : 26 Sample Multiplier: 1

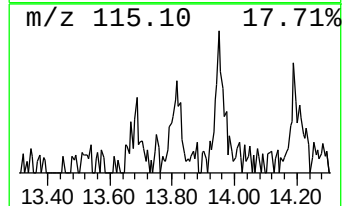
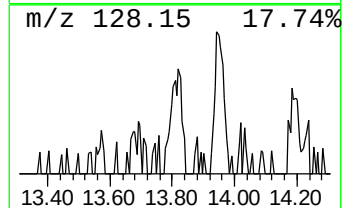
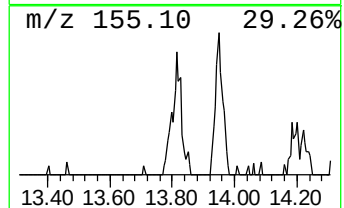
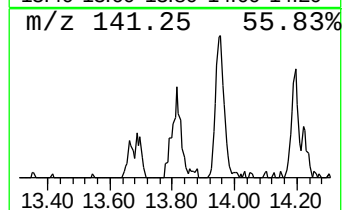
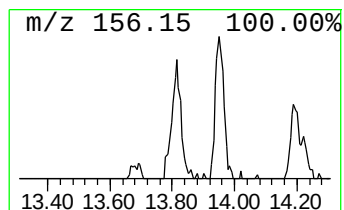
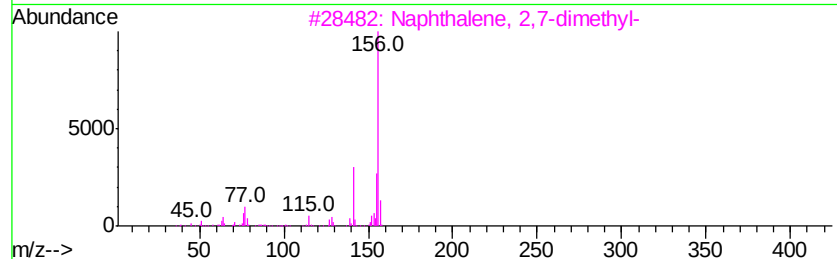
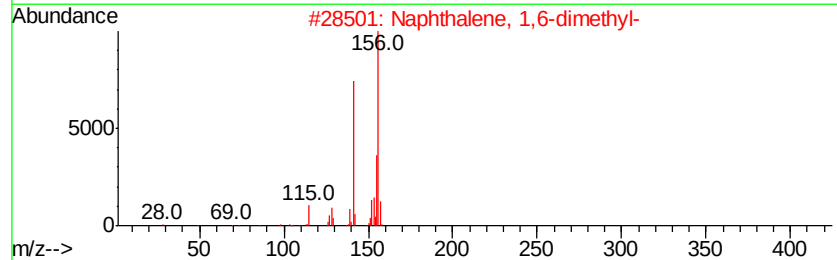
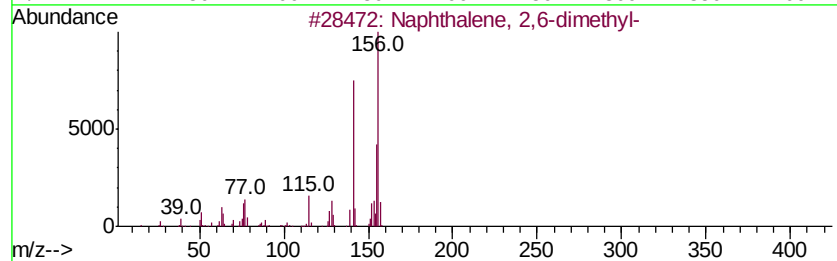
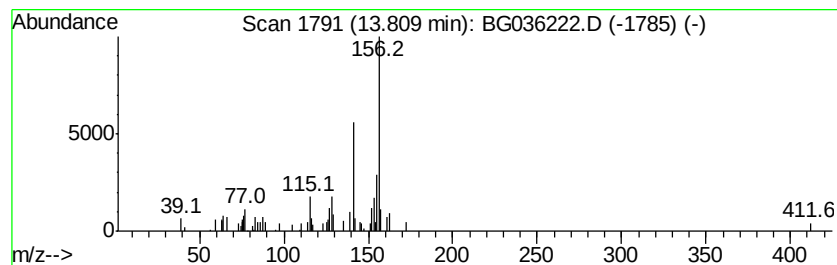
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Naphthalene, 2,6-dimethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD		R.T.
13.81	2.68 ng	49899	Acenaphthene-d10		14.63
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	94
2	Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	94
3	Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	94
4	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	93
5	Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	93



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG080818\
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Sample : J4328-01
Misc :
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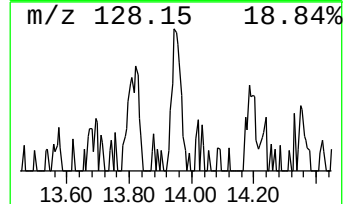
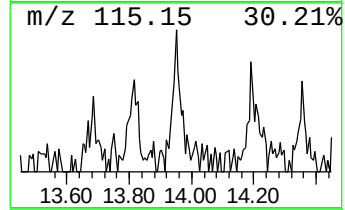
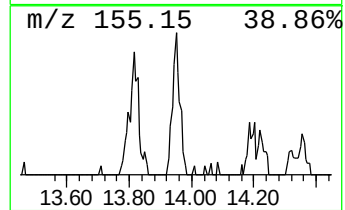
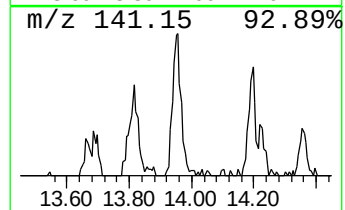
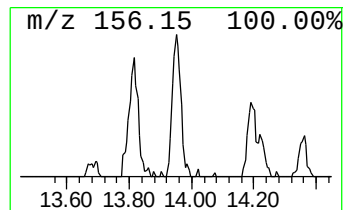
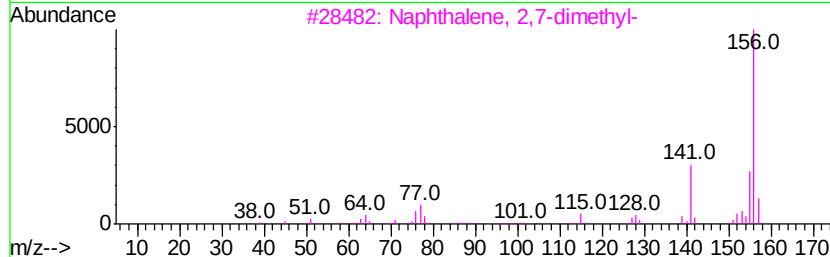
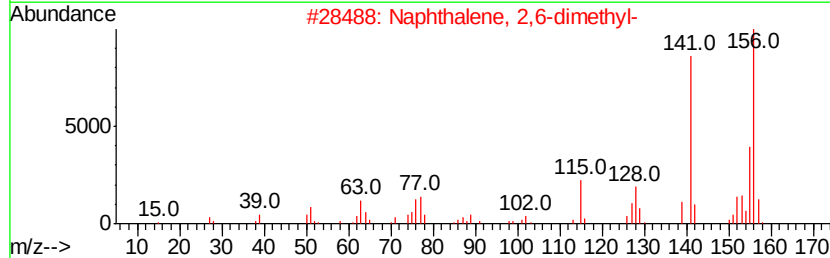
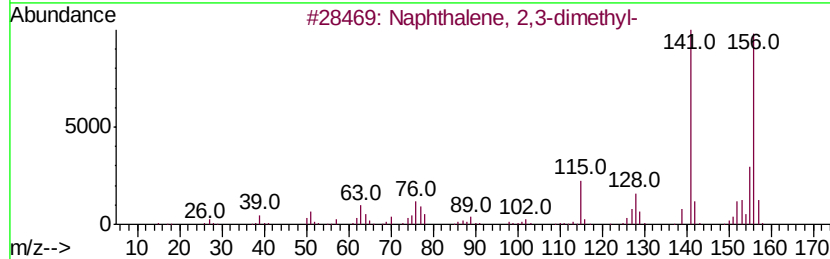
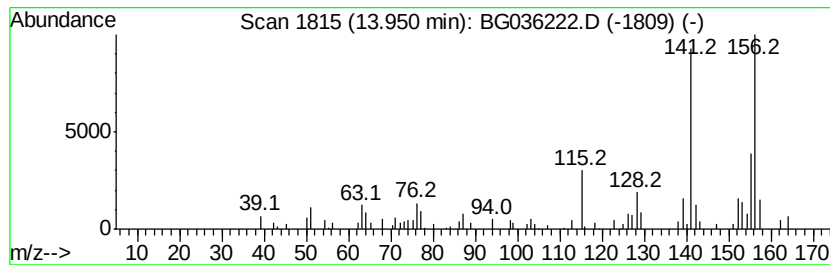
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Naphthalene, 2,3-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.95	3.46 ng	64488	Acenaphthene-d10	14.63
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 2,3-dimethyl-	156 C12H12	000581-40-8 95
2		Naphthalene, 2,6-dimethyl-	156 C12H12	000581-42-0 95
3		Naphthalene, 2,7-dimethyl-	156 C12H12	000582-16-1 95
4		Naphthalene, 1,6-dimethyl-	156 C12H12	000575-43-9 94
5		Naphthalene, 1,7-dimethyl-	156 C12H12	000575-37-1 93



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG080818\
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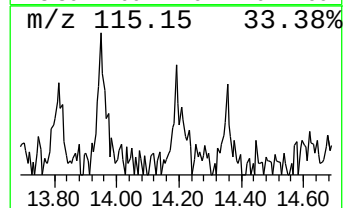
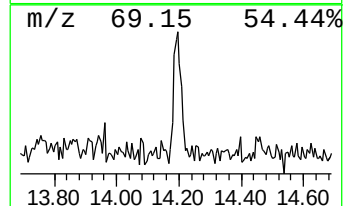
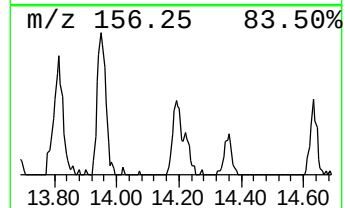
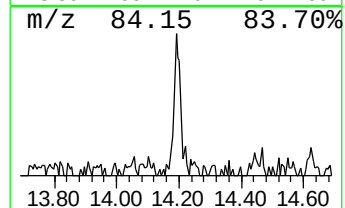
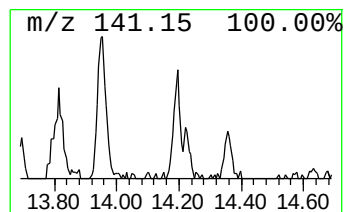
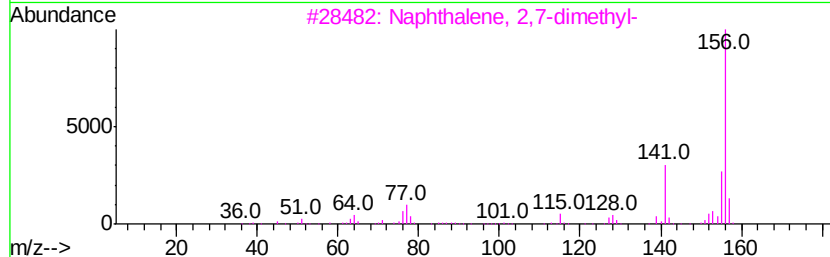
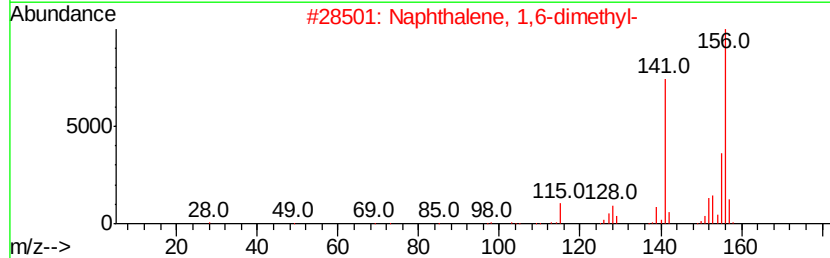
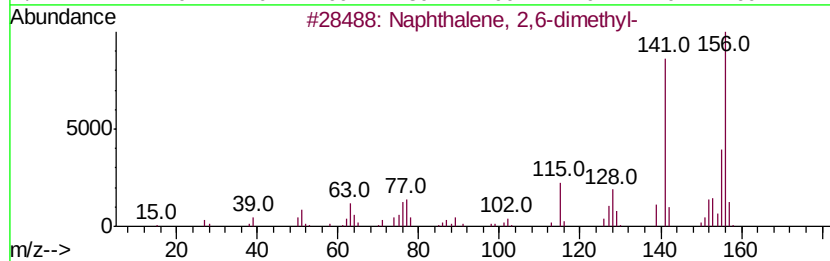
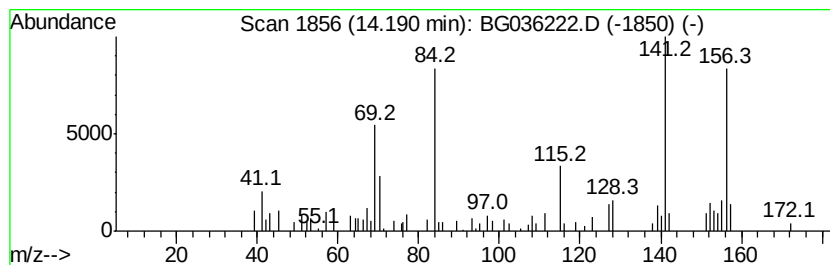
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Naphthalene, 2,7-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD		R.T.		
14.19	2.84 ng	52989	Acenaphthene-d10		14.63		
Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	78
2			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	74
3			Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	64
4			Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	64
5			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	60



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG080818\
Data File : BG036222.D
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Sample : J4328-01
Misc :
ALS Vial : 26 Sample Multiplier: 1

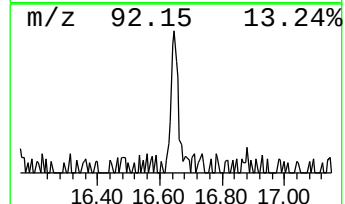
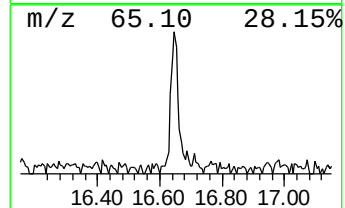
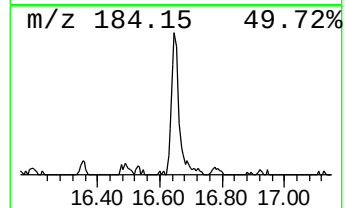
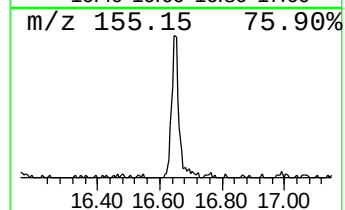
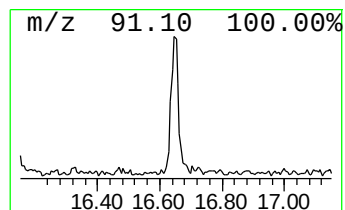
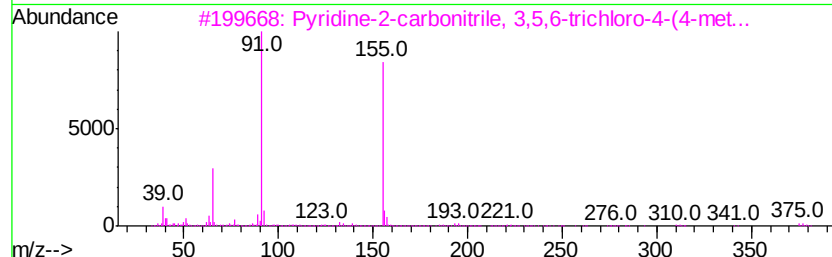
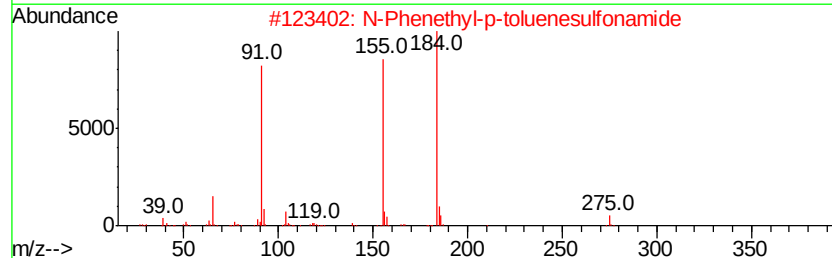
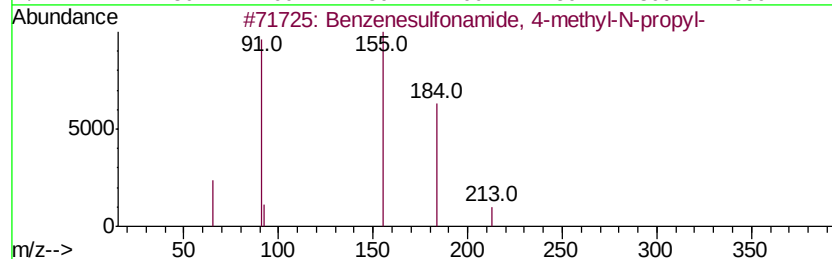
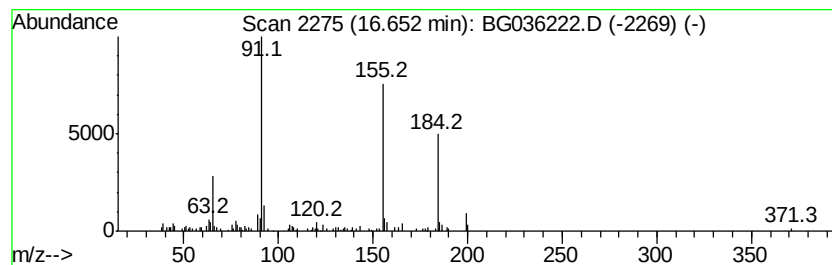
Instrument :
BNA_G
ClientSampleId :
MW-22

Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG071218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Benzenesulfonamide, 4-methy... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
16.65	4.10 ng	102335	Phenanthrene-d10	17.37	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzenesulfonamide, 4-methyl-N-p...	213	C10H15NO2S	001133-12-6	72
2	N-Phenethyl-p-toluenesulfonamide	275	C15H17NO2S	005450-75-9	64
3	Pyridine-2-carbonitrile, 3,5,6-t...	375	C13H8Cl3N3O2S	255713-77-0	58
4	(2S)-Glycidyl tosylate	228	C10H12O4S	070987-78-9	53
5	Ethyl(E/Z)-.alpha.-cyano-4-nitro...	416	C19H16N2O7S	1000287-26-8	53



Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG080818\
Data File : BG036222.D
Acq On : 9 Aug 2018 7:51
Operator : JU/SJ
Sample : J4328-01
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
MW-22

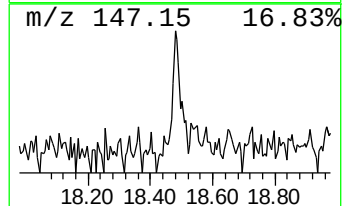
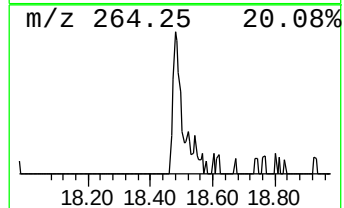
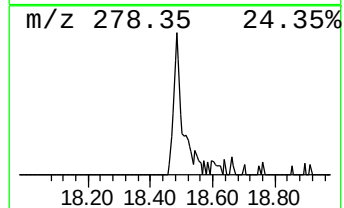
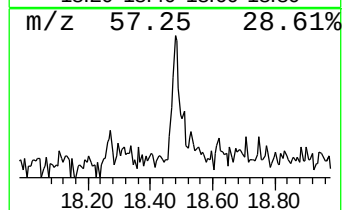
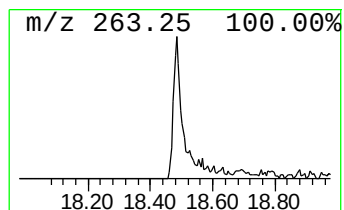
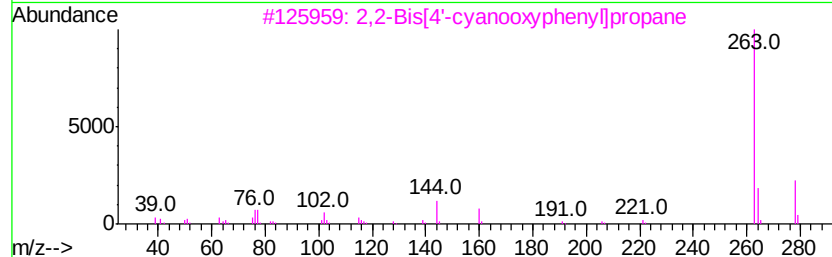
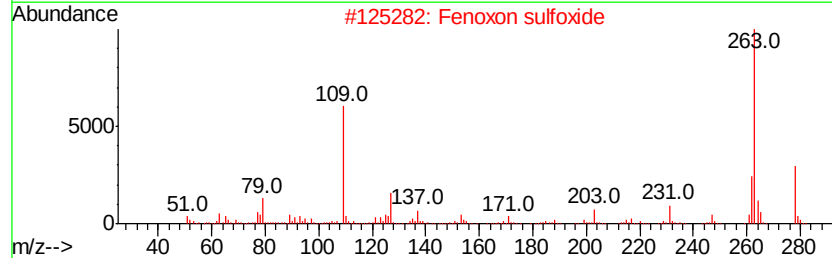
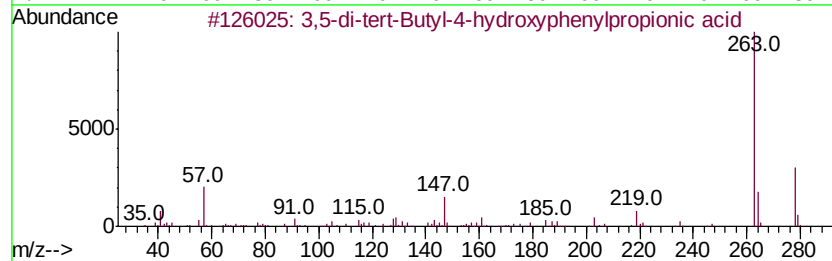
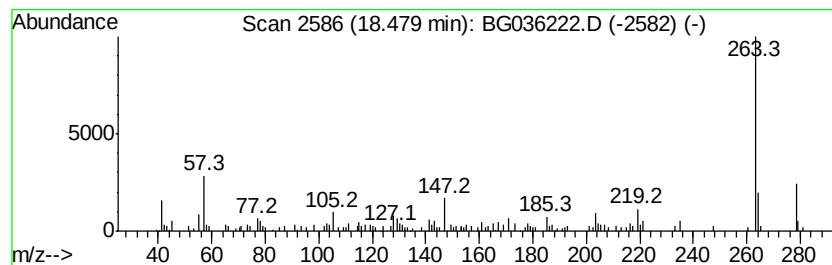
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 3,5-di-tert-Butyl-4-hydroxy... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.48	3.19 ng	79545	Phenanthrene-d10	17.37

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3,5-di-tert-Butyl-4-hydroxyphenyl...	278	C17H26O3	020170-32-5	94
2			Fenoxon sulfoxide	278	C10H15O5PS	006552-13-2	81
3			2,2-Bis[4'-cyanoxyphenyl]propane	278	C17H14N2O2	1000322-13-3	53
4			3-Phenyl-2,9-dioxo-4-aza-fluoren...	263	C16H9NO3	067138-57-2	50
5			2-(4-Methoxy-phenyl)-1H-1,3a,8-t...	263	C16H13N3O	036289-23-3	49



Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG080818\
Data File : BG036222.D
Acq On : 9 Aug 2018 7:51
Operator : JU/SJ
Sample : J4328-01
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
MW-22

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG071218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown4.64	4.64	4.4	ng	40709	1	8.01	183216	20.0
unknown5.13	5.13	4.6	ng	41809	1	8.01	183216	20.0
unknown7.55	7.55	82.5	ng	755973	1	8.01	183216	20.0
unknown9.37	9.37	3.3	ng	30154	1	8.01	183216	20.0
unknown9.51	9.51	22.4	ng	283719	2	10.81	253618	20.0
unknown9.78	9.78	2.2	ng	28018	2	10.81	253618	20.0
Benzene, (1-methy...	10.21	2.7	ng	33830	2	10.81	253618	20.0
unknown10.97	10.97	2.5	ng	31806	2	10.81	253618	20.0
Benzene, (ethenyl...	11.92	2.3	ng	29276	2	10.81	253618	20.0
Naphthalene, 2,6-...	13.81	2.7	ng	49899	3	14.63	372988	20.0
Naphthalene, 2,3-...	13.95	3.5	ng	64488	3	14.63	372988	20.0
Naphthalene, 2,7-...	14.19	2.8	ng	52989	3	14.63	372988	20.0
Benzenesulfonamid...	16.65	4.1	ng	102335	4	17.37	498620	20.0
3,5-di-tert-Butyl...	18.48	3.2	ng	79545	4	17.37	498620	20.0