

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG101015\
 Data File : BG019139.D
 Acq On : 11 Oct 2015 4:48
 Operator : UM/NP
 Sample : G3929-14
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 MW-09

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_G\METHODS\8270-BG101015.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	3.144	47	52	60	rBV	58022	97542	5.48%	1.084%
2	3.214	60	64	77	rVB	211118	282025	15.84%	3.133%
3	5.136	385	391	400	rBV	72977	104063	5.84%	1.156%
4	5.606	464	471	482	rVB	147499	231736	13.02%	2.574%
5	7.192	734	741	748	rBV	96771	160289	9.00%	1.781%
6	7.562	796	804	816	rVB	293981	490427	27.54%	5.448%
7	8.032	878	884	891	rBV	61928	100080	5.62%	1.112%
8	8.355	930	939	947	rBV	262914	448929	25.21%	4.987%
9	9.190	1073	1081	1091	rBV	257940	446017	25.05%	4.955%
10	10.835	1353	1361	1369	rVB	92500	159147	8.94%	1.768%
11	13.285	1771	1778	1790	rVB	748169	1108781	62.27%	12.318%
12	14.107	1912	1918	1927	rBV	19056	27994	1.57%	0.311%
13	14.660	2006	2012	2019	rVB2	189624	296842	16.67%	3.298%
14	16.146	2259	2265	2279	rVB	673225	1011327	56.80%	11.235%
15	16.393	2298	2307	2314	rBV3	27011	49152	2.76%	0.546%
16	16.487	2318	2323	2327	rVV	89640	121972	6.85%	1.355%
17	16.540	2327	2332	2339	rVV2	78162	161906	9.09%	1.799%
18	16.604	2339	2343	2347	rVV2	93881	129858	7.29%	1.443%
19	16.651	2347	2351	2356	rVV	52070	76138	4.28%	0.846%
20	16.698	2356	2359	2361	rVV2	16773	23836	1.34%	0.265%
21	16.728	2361	2364	2366	rVV	29757	43082	2.42%	0.479%
22	16.751	2366	2368	2372	rVV	26975	34029	1.91%	0.378%
23	16.804	2372	2377	2384	rVV4	67309	129634	7.28%	1.440%
24	16.869	2384	2388	2393	rVV	97387	140441	7.89%	1.560%
25	16.945	2397	2401	2409	rVB	45815	66516	3.74%	0.739%
26	17.404	2473	2479	2487	rVB2	231602	346803	19.48%	3.853%
27	18.073	2588	2593	2597	rBV3	16043	20177	1.13%	0.224%
28	20.018	2918	2924	2931	rVB	1378162	1780494	100.00%	19.780%
29	21.328	3143	3147	3154	rBV7	15641	32823	1.84%	0.365%
30	21.675	3200	3206	3213	rBV	262344	429719	24.13%	4.774%
31	24.901	3745	3755	3768	rVB	145965	405890	22.80%	4.509%
32	25.071	3778	3784	3795	rVB5	15957	43688	2.45%	0.485%

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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_G\METHODS\8270-BG101015.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

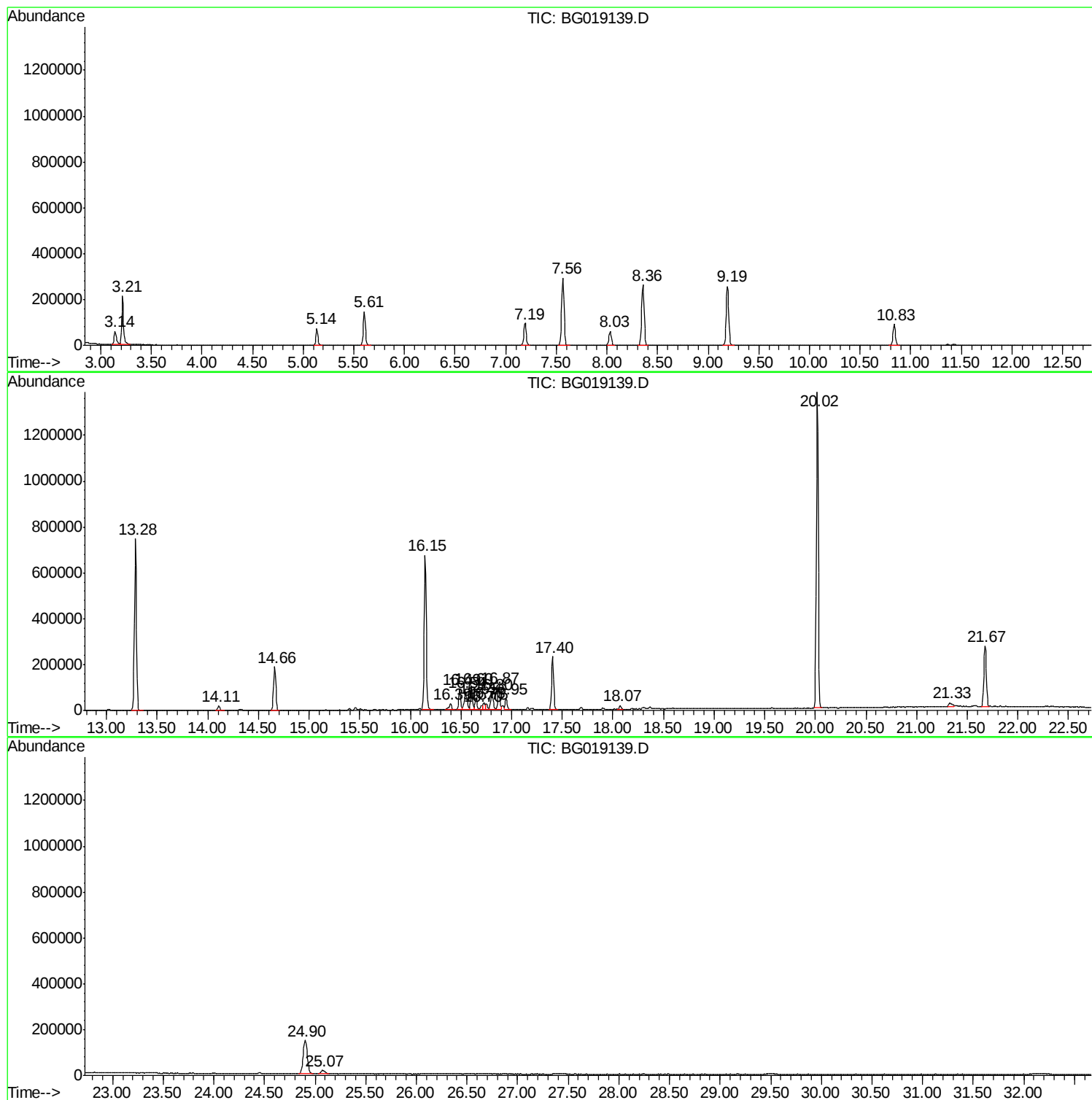
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Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG101015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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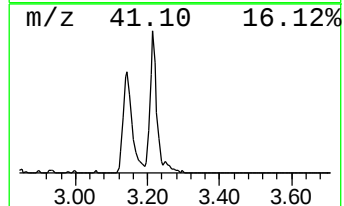
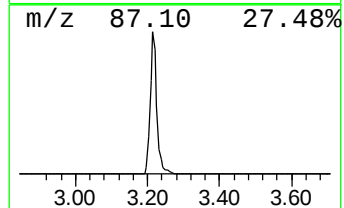
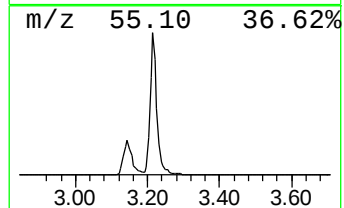
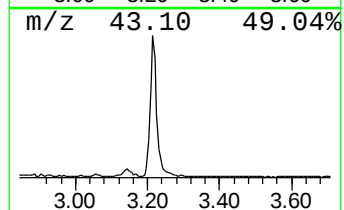
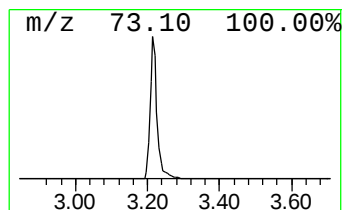
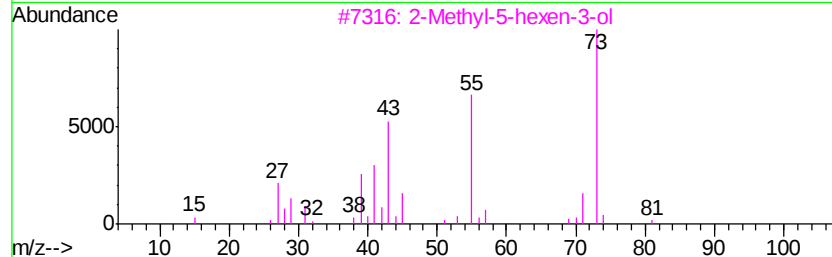
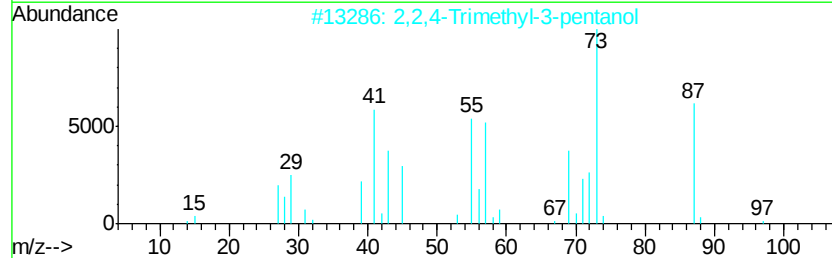
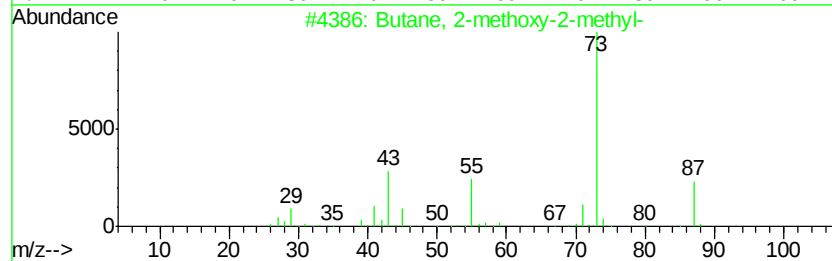
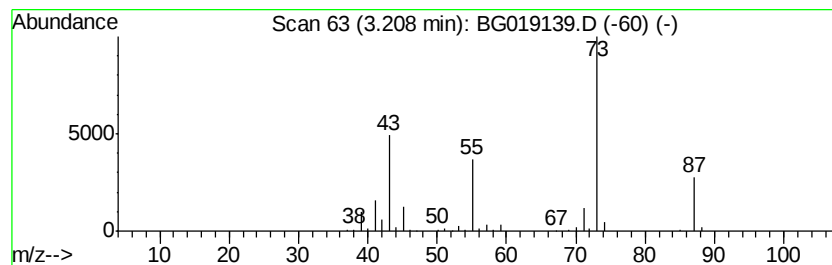
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Butane, 2-methoxy-2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.21	56.36 ng	282025	1,4-Dichlorobenzene-d4	8.03

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
3		2-Methyl-5-hexen-3-ol	114	C7H14O	032815-70-6	37
4		1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	9
5		3-Pentanol, 2,4-dimethyl-	116	C7H16O	000600-36-2	9



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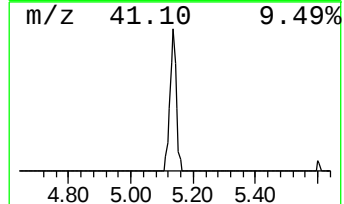
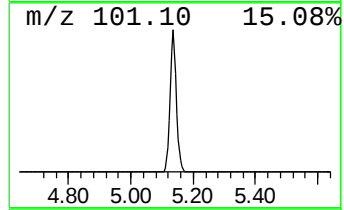
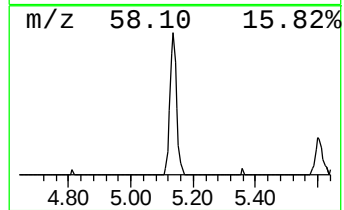
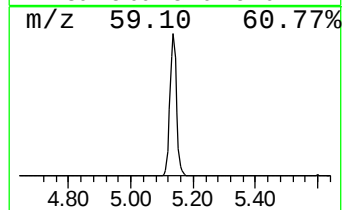
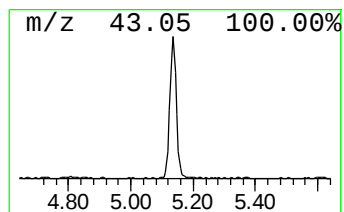
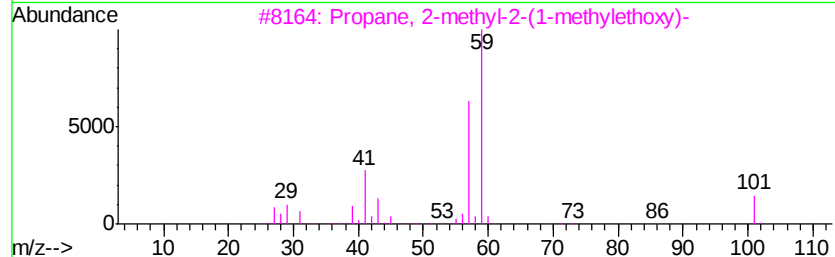
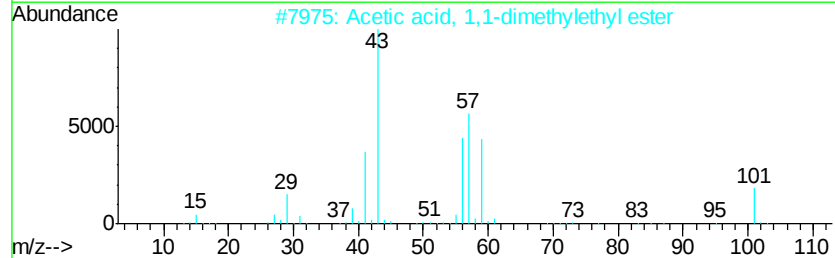
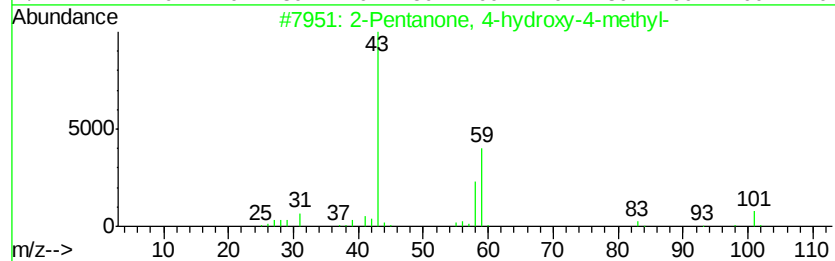
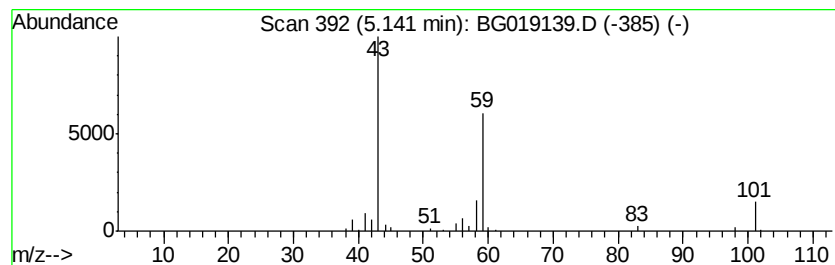
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.14	20.80 ng	104063	1,4-Dichlorobenzene-d4	8.03

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2		000123-42-2	50
2	Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2		000540-88-5	38
3	Propane, 2-methyl-2-(1-methyleth...	116	C7H16O		017348-59-3	36
4	3-Hexanol, 4-methyl-	116	C7H16O		000615-29-2	33
5	2-Hexanol, 2-methyl-	116	C7H16O		000625-23-0	28



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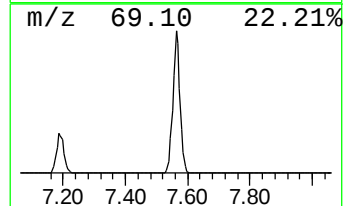
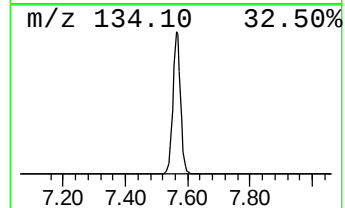
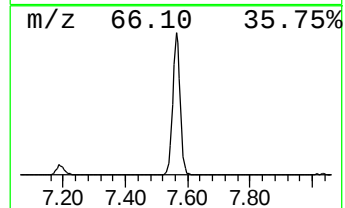
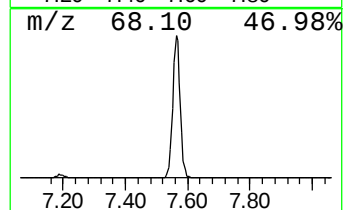
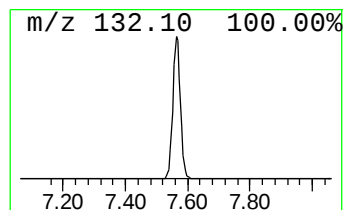
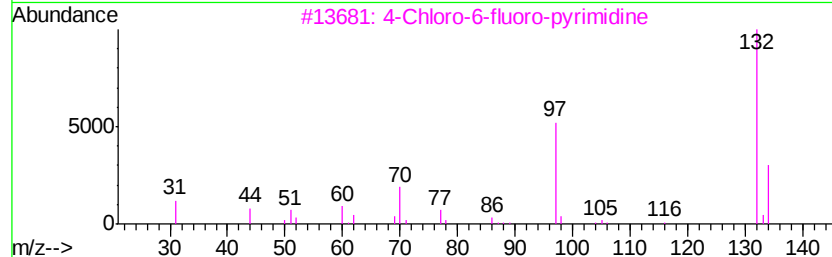
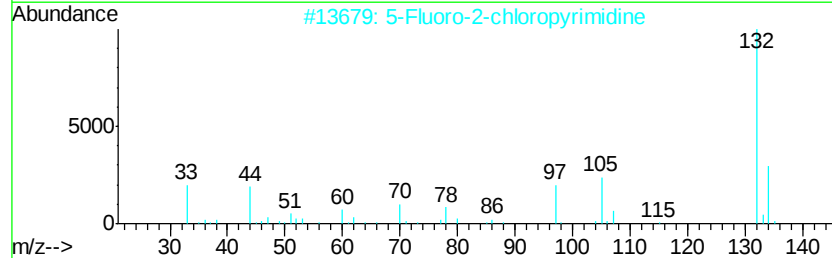
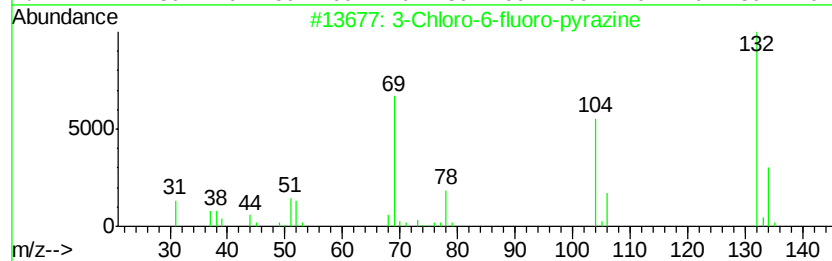
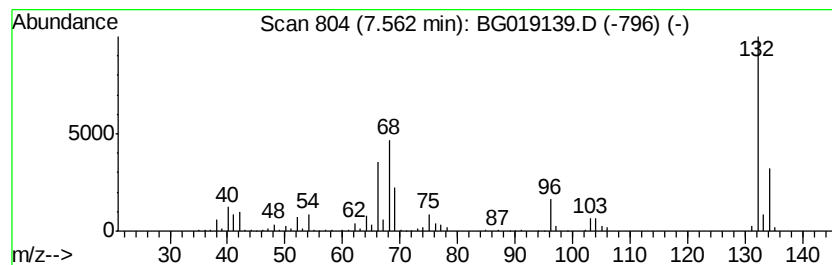
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 unknown7.56 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.56	98.01 ng	490427	1,4-Dichlorobenzene-d4	8.03

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	27
3		4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	27
4		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
5		1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	17



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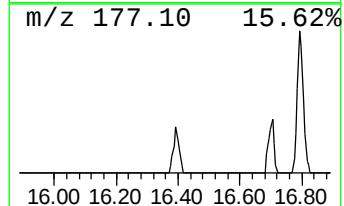
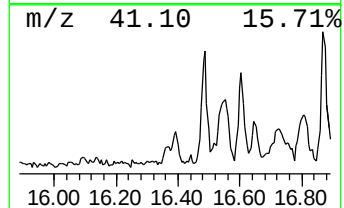
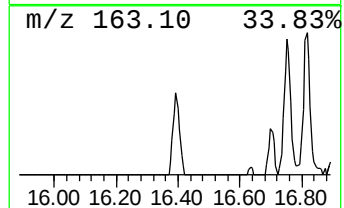
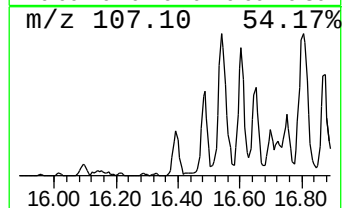
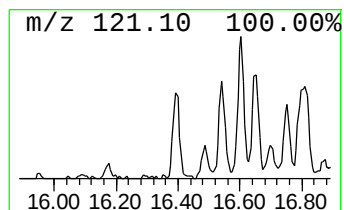
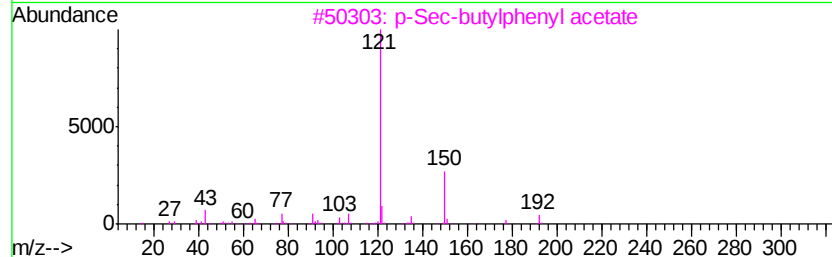
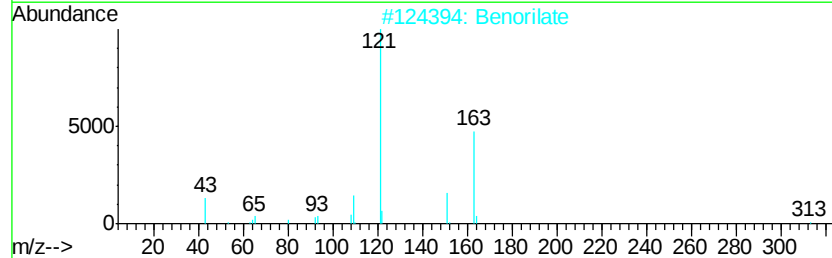
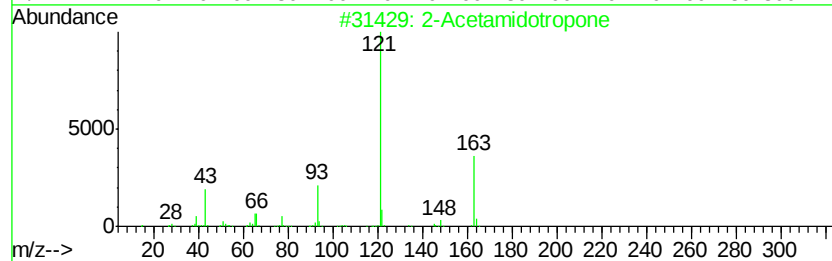
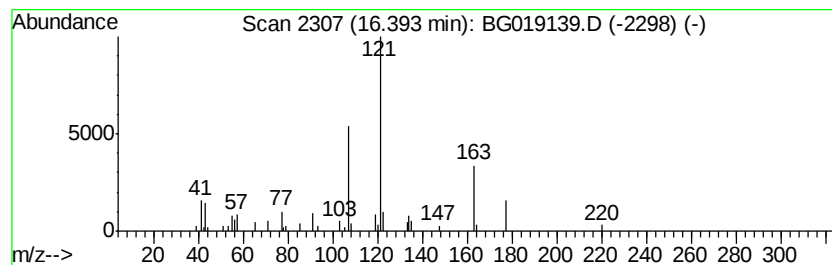
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 unknown16.39 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD		R.T.
16.39	2.83 ng	49152	Phenanthrene-d10		17.40

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2		Acetamidotropone	163	C9H9NO2	006422-12-4	47
2			Benorilate	313	C17H15NO5	005003-48-5	47
3			p-Sec-butylphenyl acetate	192	C12H16O2	003245-24-7	43
4			3',4'-Acetoxylide	163	C10H13NO	002198-54-1	40
5			Benzenamine, 2,5-dimethyl-	121	C8H11N	000095-78-3	38



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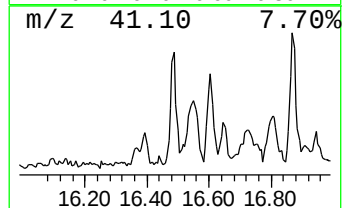
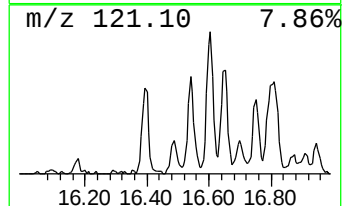
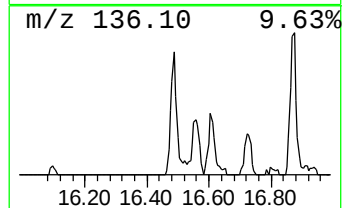
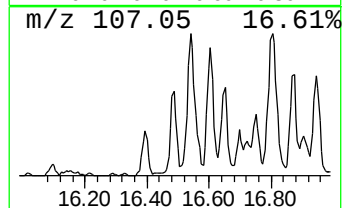
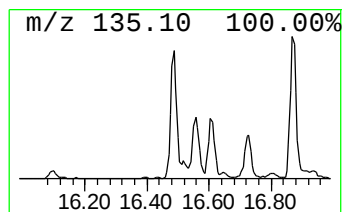
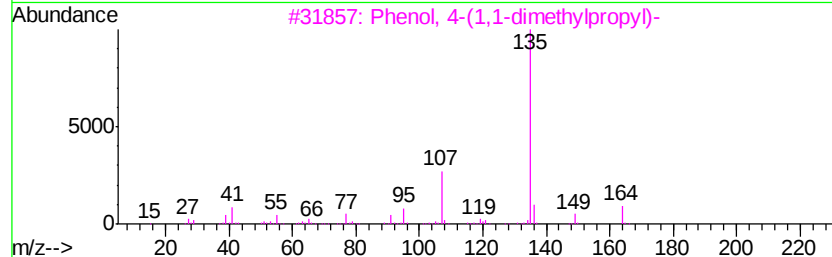
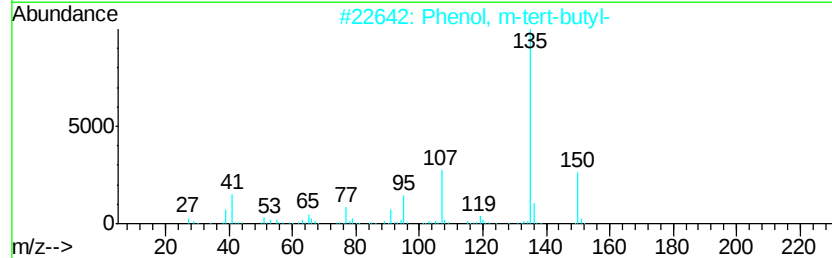
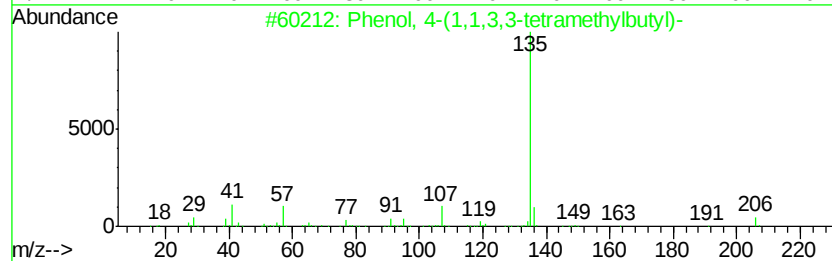
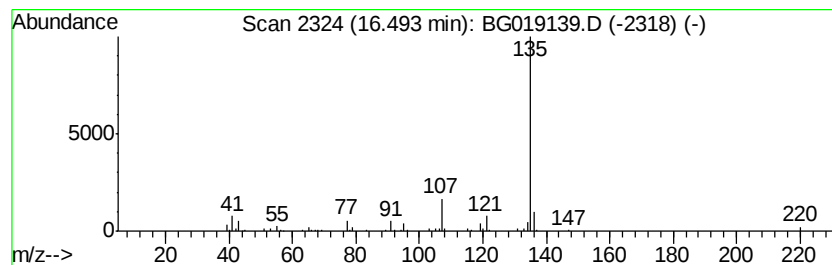
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 7 Phenol, 4-(1,1,3,3-tetramet... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.49	7.03 ng	121972	Phenanthrene-d10	17.40

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phenol, 4-(1,1,3,3-tetramethylbu...	206	C14H22O	000140-66-9	78
2		Phenol, m-tert-butyl-	150	C10H14O	000585-34-2	78
3		Phenol, 4-(1,1-dimethylpropyl)-	164	C11H16O	000080-46-6	78
4		Hexestrol	270	C18H22O2	005635-50-7	64
5		1-n-Hexyladamantane	220	C16H28	022458-75-9	59



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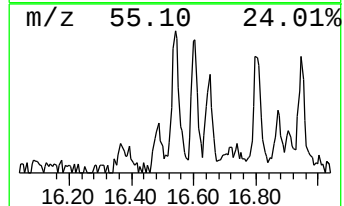
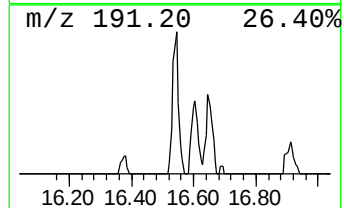
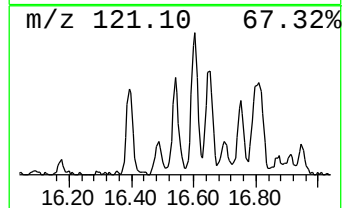
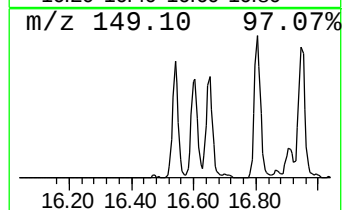
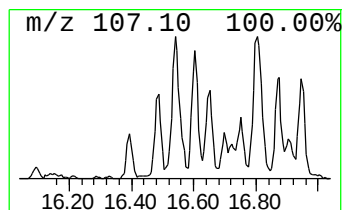
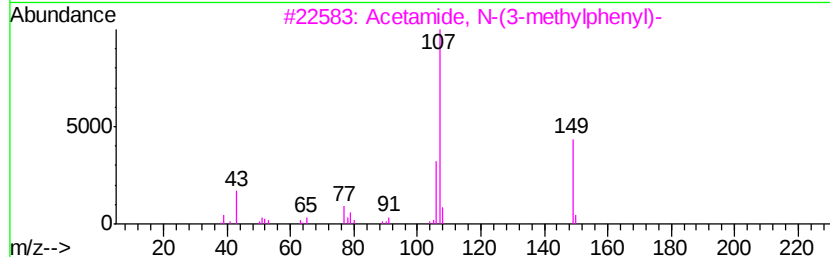
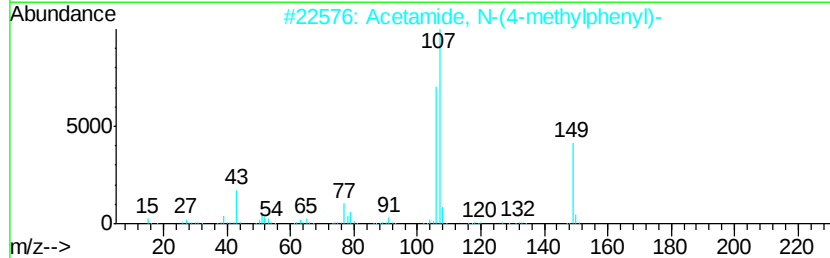
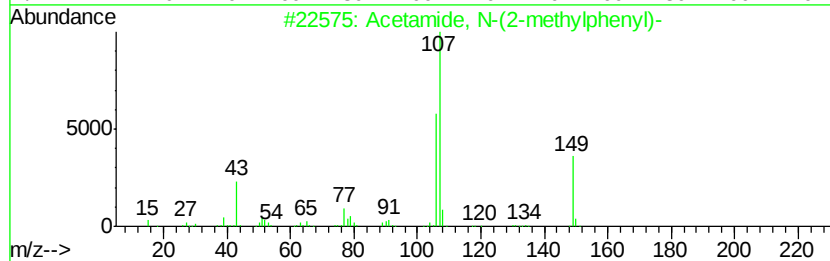
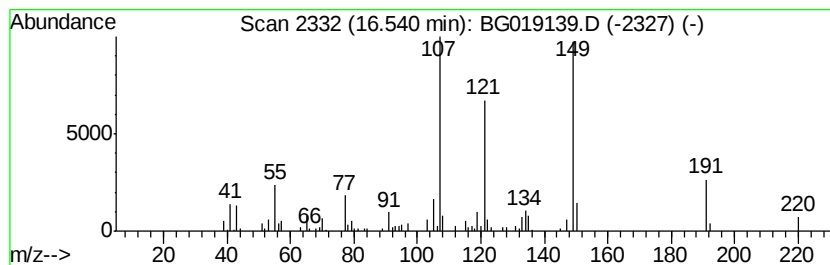
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 unknown16.54 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.54	9.34 ng	161906	Phenanthrene-d10	17.40

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Acetamide, N-(2-methylphenyl)-	149	C9H11NO	000120-66-1	46
2			Acetamide, N-(4-methylphenyl)-	149	C9H11NO	000103-89-9	43
3			Acetamide, N-(3-methylphenyl)-	149	C9H11NO	000537-92-8	42
4			Ethanone, 2-(1-methylethoxy)-1,2...	254	C17H18O2	006652-28-4	38
5			4-Methyl-2-tert-octylphenol	220	C15H24O	004979-46-8	38



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG101015\
Data File : BG019139.D
Acq On : 11 Oct 2015 4:48
Operator : UM/NP
Sample : G3929-14
Misc :
ALS Vial : 24 Sample Multiplier: 1

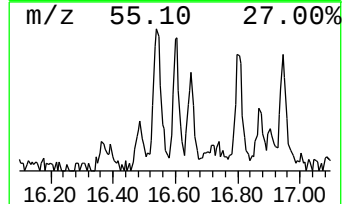
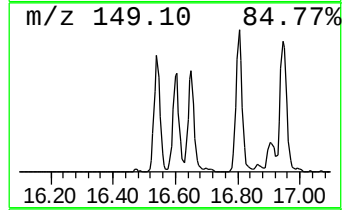
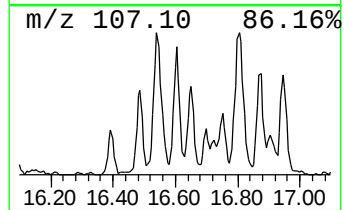
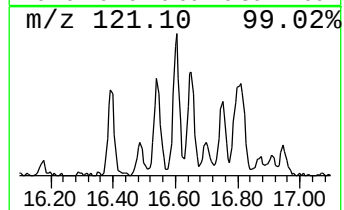
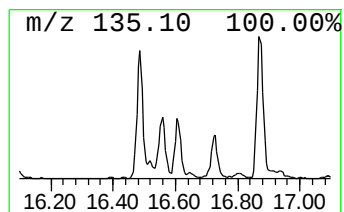
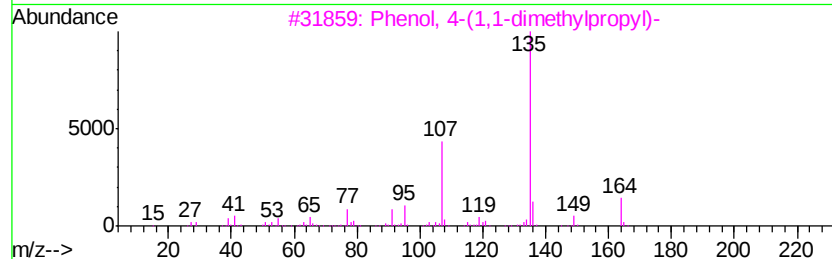
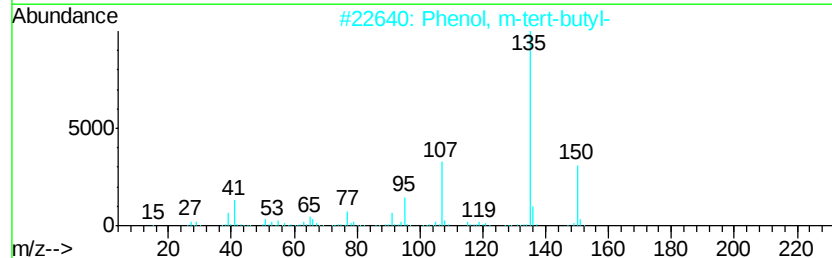
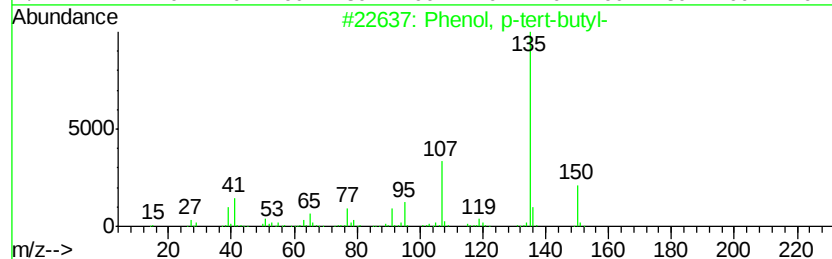
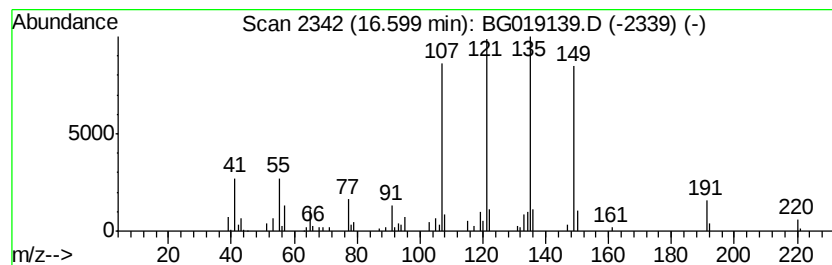
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ClientSampled :
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Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG101015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Phenol, p-tert-butyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD		R.T.	
16.60	7.49 ng	129858	Phenanthrene-d10		17.40	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	64	
2	Phenol, m-tert-butyl-	150	C10H14O	000585-34-2	52	
3	Phenol, 4-(1,1-dimethylpropyl)-	164	C11H16O	000080-46-6	50	
4	Phenol, 4,4'-(1,2-diethyl-1,2-et...	270	C18H22O2	000084-16-2	47	
5	Benzeneacetonitrile, 4-fluoro-	135	C8H6FN	000459-22-3	43	



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG101015\
Data File : BG019139.D
Acq On : 11 Oct 2015 4:48
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Sample : G3929-14
Misc :
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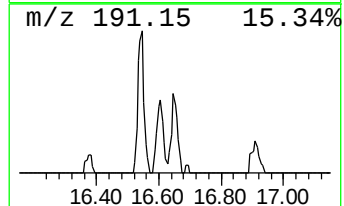
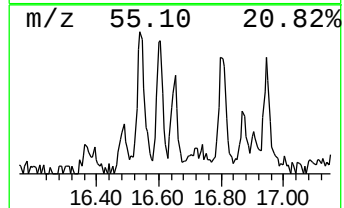
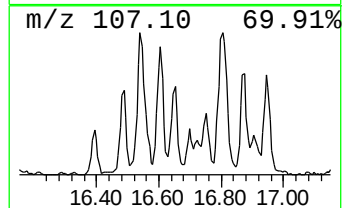
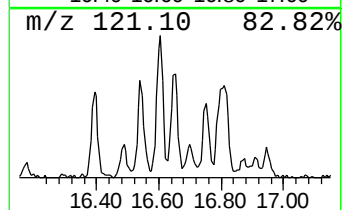
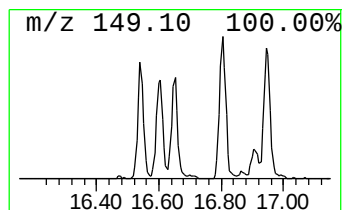
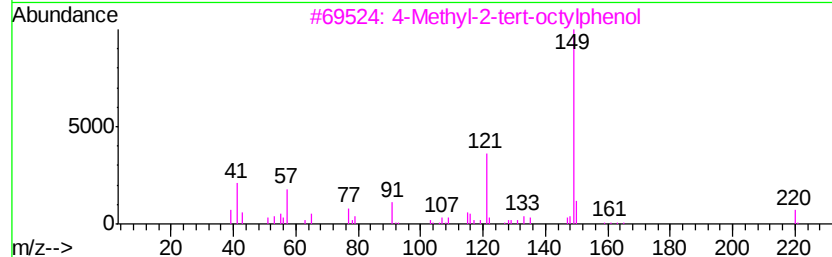
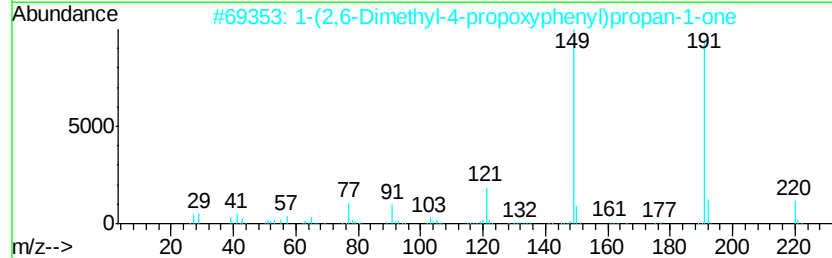
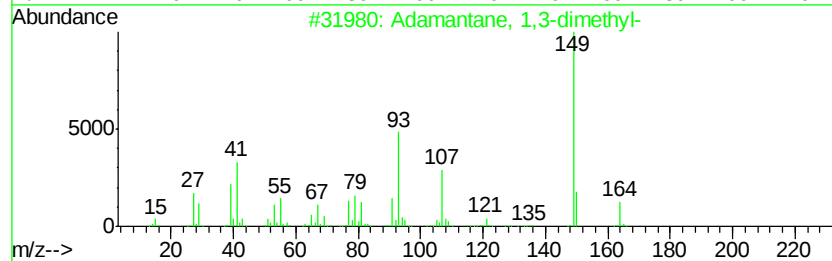
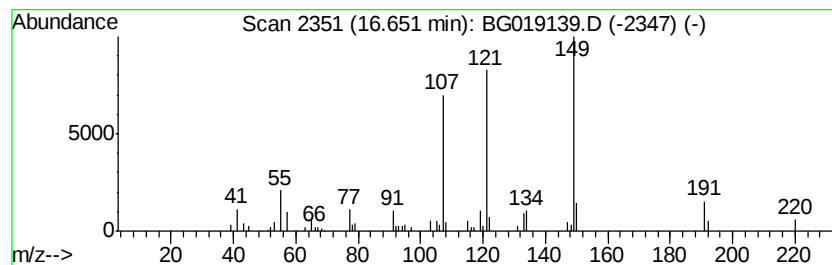
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG101015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 unknown16.65 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.65	4.39 ng	76138	Phenanthrene-d10	17.40
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Adamantane, 1,3-dimethyl-	164 C12H20	000702-79-4 35
2		1-(2,6-Dimethyl-4-propoxyphenyl)...	220 C14H20O2	1000190-22-3 35
3		4-Methyl-2-tert-octylphenol	220 C15H24O	004979-46-8 30
4		2-Methyl-2-(4'-methoxyphenyl)pen...	206 C13H18O2	185700-49-6 27
5		Phenol, 2-(1-methylpropyl)-	150 C10H14O	000089-72-5 27



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG101015\
Data File : BG019139.D
Acq On : 11 Oct 2015 4:48
Operator : UM/NP
Sample : G3929-14
Misc :
ALS Vial : 24 Sample Multiplier: 1

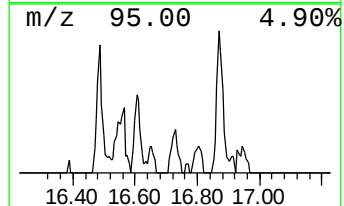
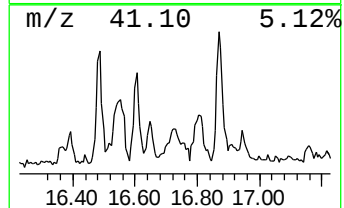
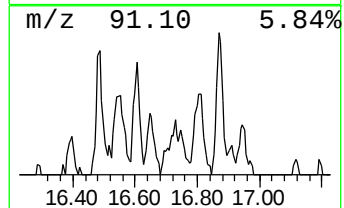
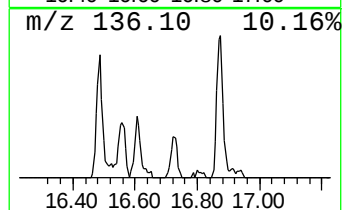
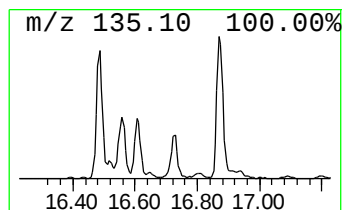
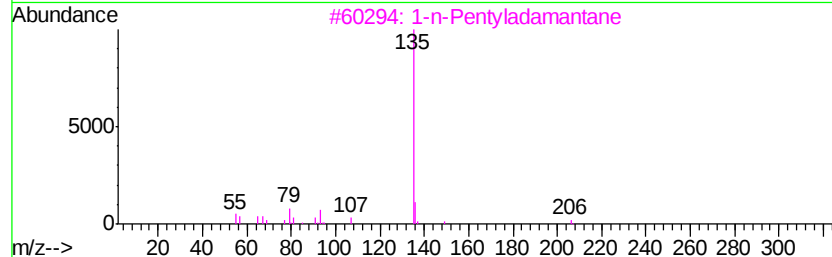
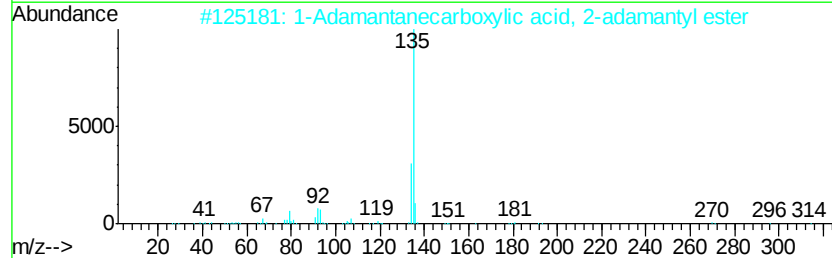
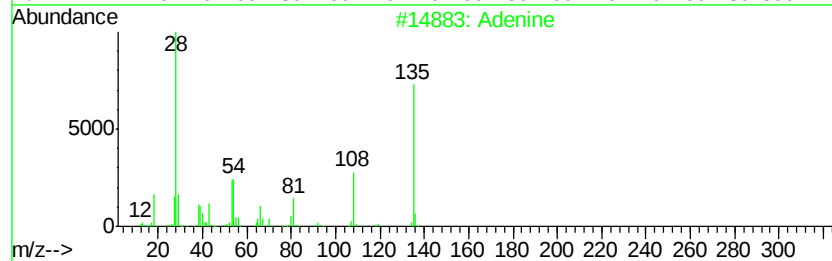
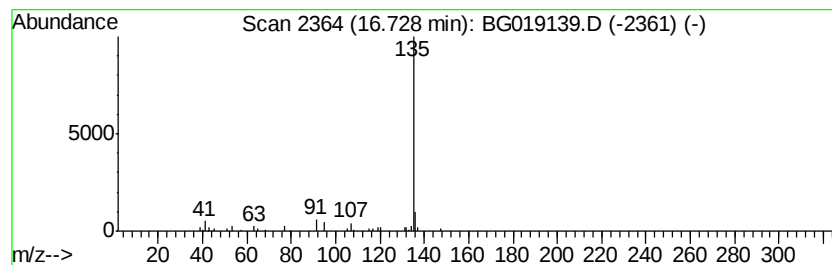
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Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG101015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Adenine Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD		R.T.	
16.73	2.48 ng	43082	Phenanthrene-d10		17.40	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Adenine		135	C5H5N5	000073-24-5	56
2	1-Adamantanecarboxylic acid, 2-a...		314	C21H30O2	1000282-94-3	50
3	1-n-Pentyladamantane		206	C15H26	050782-11-1	40
4	Silane, trichloro(2-tricyclo[3.3...		296	C12H19Cl3Si	037843-11-1	40
5	Butanenitrile, 2-(1-adamantyl)-3...		217	C15H23N	1000266-44-6	40



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG101015\
Data File : BG019139.D
Acq On : 11 Oct 2015 4:48
Operator : UM/NP
Sample : G3929-14
Misc :
ALS Vial : 24 Sample Multiplier: 1

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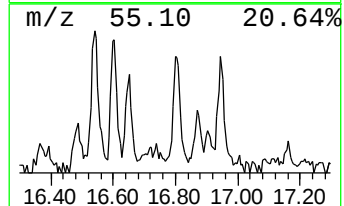
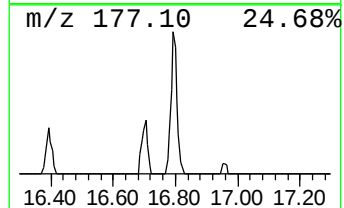
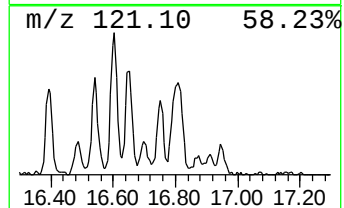
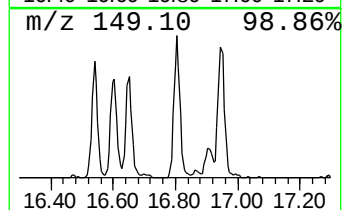
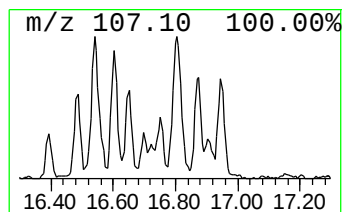
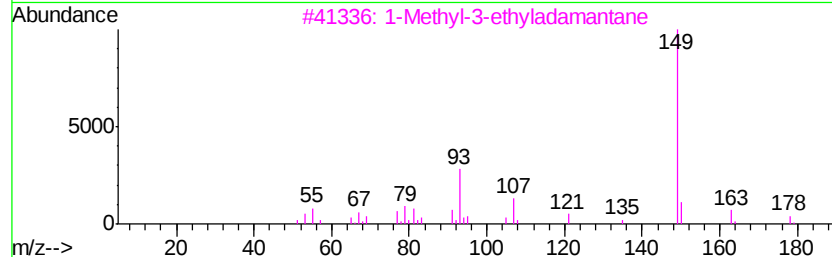
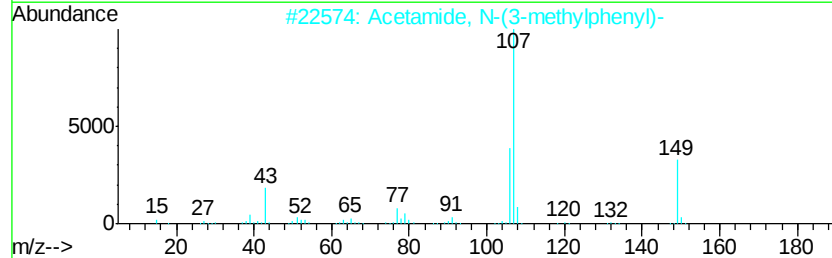
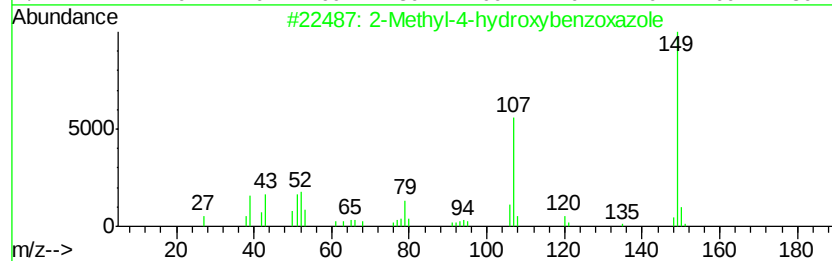
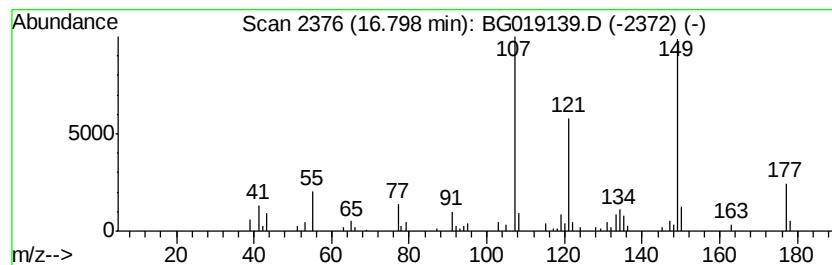
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 2-Methyl-4-hydroxybenzoxazole Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.80	7.48 ng	129634	Phenanthrene-d10	17.40

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Methyl-4-hydroxybenzoxazole	149	C8H7NO2	051110-60-2	59
2			Acetamide, N-(3-methylphenyl)-	149	C9H11NO	000537-92-8	58
3			1-Methyl-3-ethyladamantane	178	C13H22	001687-34-9	43
4			N,N,2,4-Tetramethylaniline #	149	C10H15N	000769-53-9	38
5			Phenol, 2-methyl-4-(1,1,3,3-tetr...	220	C15H24O	002219-84-3	38



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG101015\
Data File : BG019139.D
Acq On : 11 Oct 2015 4:48
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Sample : G3929-14
Misc :
ALS Vial : 24 Sample Multiplier: 1

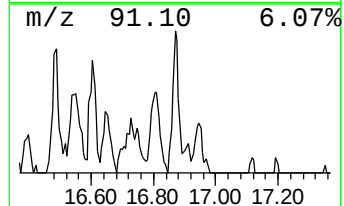
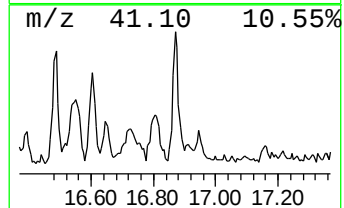
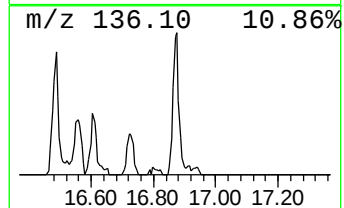
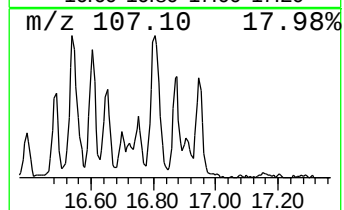
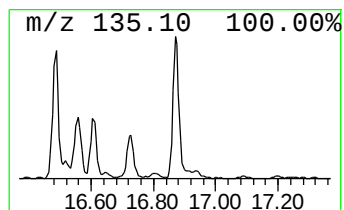
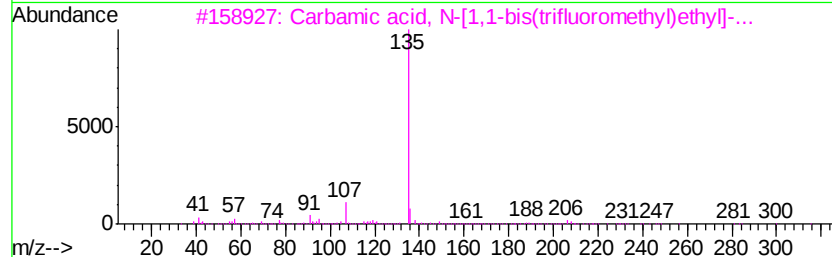
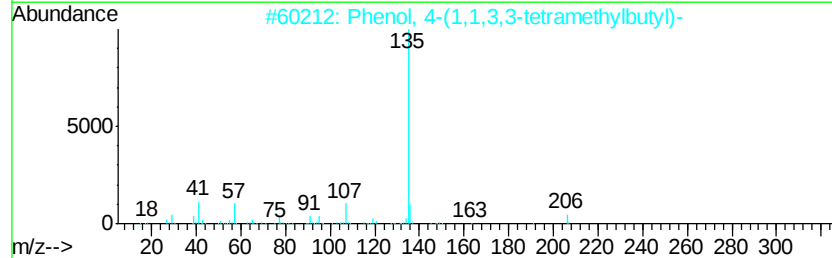
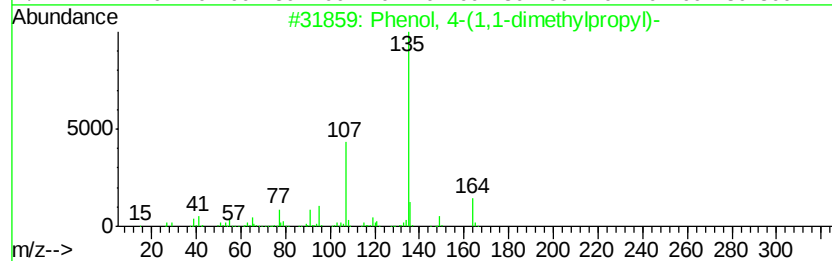
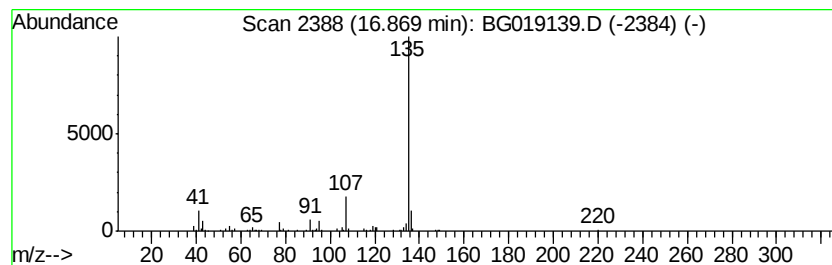
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG101015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Phenol, 4-(1,1-dimethylprop... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.		
16.87	8.10 ng	140441	Phenanthrene-d10	17.40		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 4-(1,1-dimethylpropyl)-	164	C11H16O	000080-46-6	78	
2	Phenol, 4-(1,1,3,3-tetramethylbu...	206	C14H22O	000140-66-9	78	
3	Carbamic acid, N-[1,1-bis(triflu...	413	C19H25F6N02	296242-69-8	72	
4	Phenol, 2-methyl-5-(1-methylethyl)-	150	C10H14O	000499-75-2	64	
5	Benzoic acid, 4-(bromomethyl)-	214	C8H7BrO2	006232-88-8	64	



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG101015\
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Acq On : 11 Oct 2015 4:48
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Misc :
ALS Vial : 24 Sample Multiplier: 1

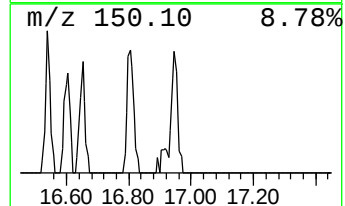
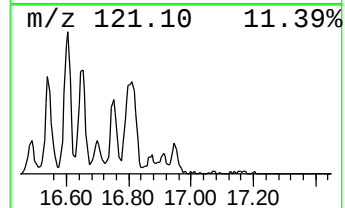
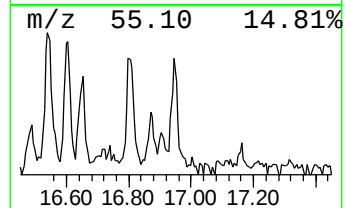
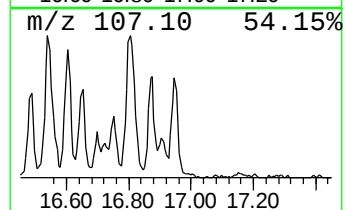
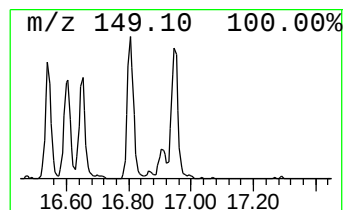
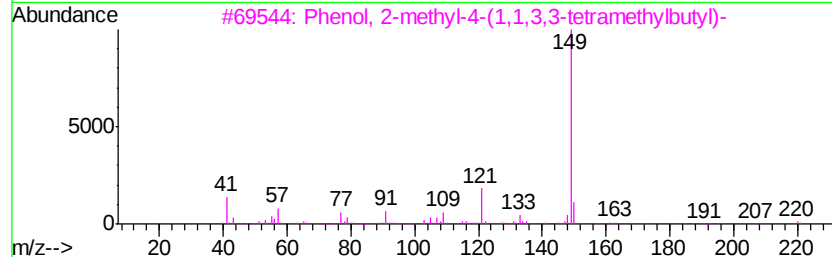
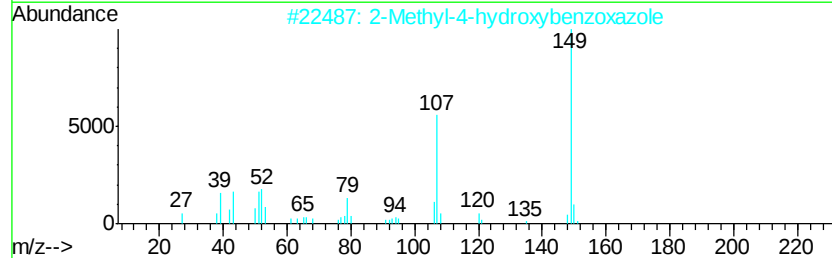
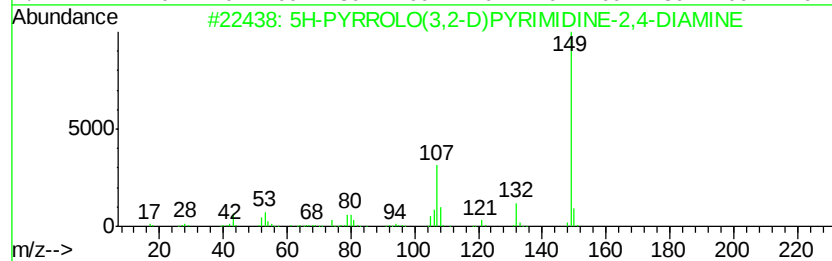
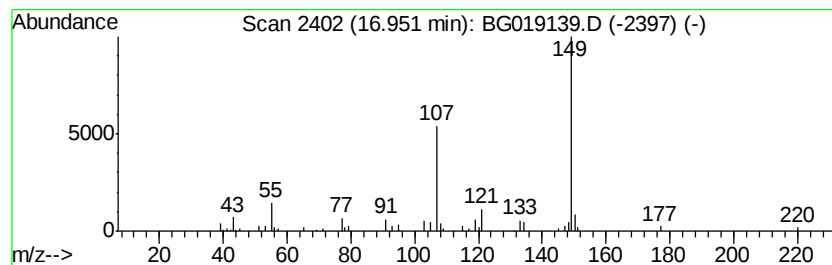
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG101015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 5H-PYRROLO(3,2-D)PYRIMIDINE... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.		
16.95	3.84 ng	66516	Phenanthrene-d10	17.40		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	5H-PYRROLO(3,2-D)PYRIMIDINE-2,4-...	149	C6H7N5	1000244-21-4	72	
2	2-Methyl-4-hydroxybenzoxazole	149	C8H7N02	051110-60-2	64	
3	Phenol, 2-methyl-4-(1,1,3,3-tetr...	220	C15H24O	002219-84-3	53	
4	Pyridine-3-carboxamide, 4-dimeth...	277	C14H13F2N3O	1000266-41-1	40	
5	Benzo[1,3]dioxole-5-carboxylic a...	362	C16H14N2O6S	1000296-89-7	38	



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG101015\
Data File : BG019139.D
Acq On : 11 Oct 2015 4:48
Operator : UM/NP
Sample : G3929-14
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
MW-09

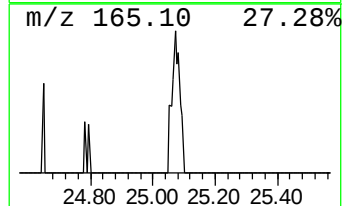
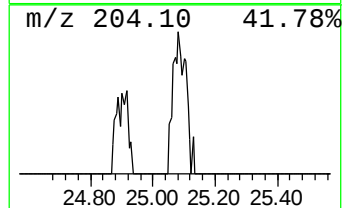
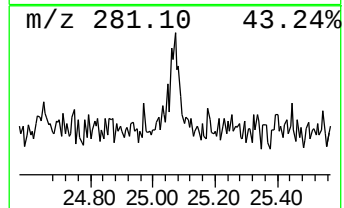
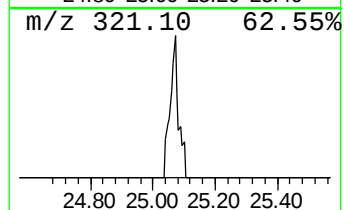
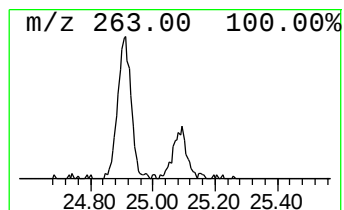
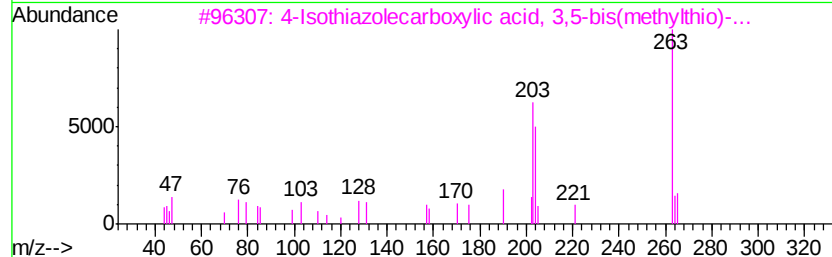
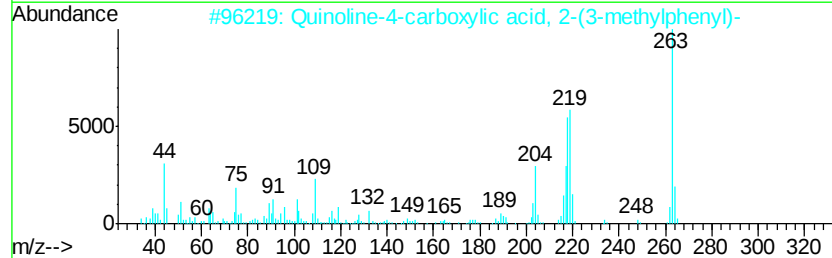
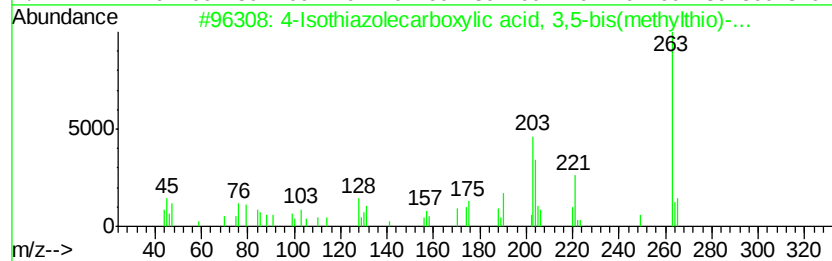
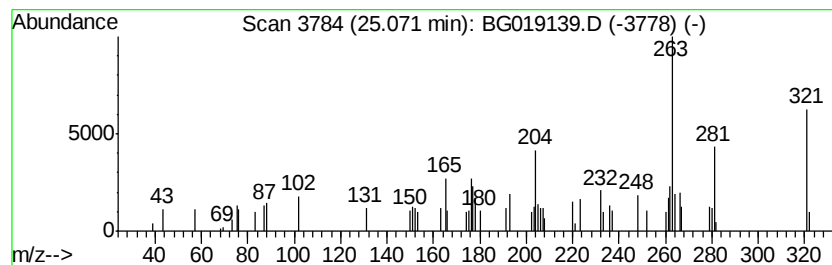
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 unknown25.07 Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.07	2.15 ng	43688	Perylene-d12	24.90

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4-Isothiazolecarboxylic acid, 3,...	263	C9H13N02S3	037572-39-7	38
2		Quinoline-4-carboxylic acid, 2-(...	263	C17H13N02	020389-04-2	30
3		4-Isothiazolecarboxylic acid, 3,...	263	C9H13N02S3	037572-40-0	30
4		1H-Indole, 2-(2'-tetrahydrofury...	263	C18H17N0	163064-85-5	30
5		3,4-Dihydrocoumarin, 4,4,5,7,8-p...	263	C14H17N04	040662-17-7	27



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG101015\
Data File : BG019139.D
Acq On : 11 Oct 2015 4:48
Operator : UM/NP
Sample : G3929-14
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
MW-09

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG101015.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...	3.21	56.4	ng	282025	1	8.03	100080	20.0
2-Pentanone, 4-hy...	5.14	20.8	ng	104063	1	8.03	100080	20.0
unknown7.56	7.56	98.0	ng	490427	1	8.03	100080	20.0
unknown16.39	16.39	2.8	ng	49152	4	17.40	346803	20.0
Phenol, 4-(1,1,3,...	16.49	7.0	ng	121972	4	17.40	346803	20.0
unknown16.54	16.54	9.3	ng	161906	4	17.40	346803	20.0
Phenol, p-tert-bu...	16.60	7.5	ng	129858	4	17.40	346803	20.0
unknown16.65	16.65	4.4	ng	76138	4	17.40	346803	20.0
Adenine	16.73	2.5	ng	43082	4	17.40	346803	20.0
2-Methyl-4-hydrox...	16.80	7.5	ng	129634	4	17.40	346803	20.0
Phenol, 4-(1,1-di...	16.87	8.1	ng	140441	4	17.40	346803	20.0
5H-PYRROLO(3,2-D)...	16.95	3.8	ng	66516	4	17.40	346803	20.0
unknown25.07	25.07	2.1	ng	43688	6	24.90	405890	20.0