

Data Path : Z:\HPCHEM1\BNA G\DATA\BG122216\
 Data File : BG025386.D
 Acq On : 22 Dec 2016 12:48
 Operator : UM/SJ
 Sample : H6199-02
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 W-05

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-BG122116.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	5.271	409	414	427	rVB	114163	193194	3.28%	0.780%
2	5.753	489	496	507	rBV	394015	712551	12.08%	2.878%
3	7.374	765	772	785	rBV	261602	511786	8.68%	2.067%
4	7.744	826	835	847	rBV	788685	1453682	24.64%	5.872%
5	8.214	908	915	921	rBV2	188901	344696	5.84%	1.392%
6	8.538	959	970	978	rVB	693145	1274385	21.60%	5.147%
7	9.395	1107	1116	1126	rBV	672012	1288745	21.85%	5.205%
8	9.513	1128	1136	1145	rVB6	47977	112022	1.90%	0.452%
9	10.206	1247	1254	1260	rBV3	87620	183969	3.12%	0.743%
10	11.041	1388	1396	1407	rBV	263088	505099	8.56%	2.040%
11	12.022	1556	1563	1572	rBV2	96598	184077	3.12%	0.744%
12	13.461	1799	1808	1814	rBV	1837799	2988117	50.65%	12.070%
13	14.184	1923	1931	1941	rBV6	78107	194456	3.30%	0.785%
14	14.278	1941	1947	1951	rVV2	123294	185157	3.14%	0.748%
15	14.836	2034	2042	2049	rBV	475984	804666	13.64%	3.250%
16	15.277	2113	2117	2124	rBV8	42977	75122	1.27%	0.303%
17	15.441	2136	2145	2151	rVB8	52060	135612	2.30%	0.548%
18	15.576	2164	2168	2172	rVB4	77051	112683	1.91%	0.455%
19	16.182	2266	2271	2279	rVB6	64343	123651	2.10%	0.499%
20	16.328	2288	2296	2307	rVB	2050382	3131676	53.09%	12.649%
21	16.493	2319	2324	2331	rBV10	77345	143581	2.43%	0.580%
22	17.586	2504	2510	2518	rVB	593996	934989	15.85%	3.777%
23	17.915	2562	2566	2574	rVB3	72767	129293	2.19%	0.522%
24	18.420	2648	2652	2657	rBV5	80059	120399	2.04%	0.486%
25	19.677	2863	2866	2869	rBV4	46259	60582	1.03%	0.245%
26	19.795	2883	2886	2892	rVB2	53354	98898	1.68%	0.399%
27	20.165	2943	2949	2955	rVB	4524236	5898966	100.00%	23.827%
28	21.875	3234	3240	3249	rVB	829372	1431109	24.26%	5.781%
29	25.283	3810	3820	3834	rVB2	467975	1424362	24.15%	5.753%

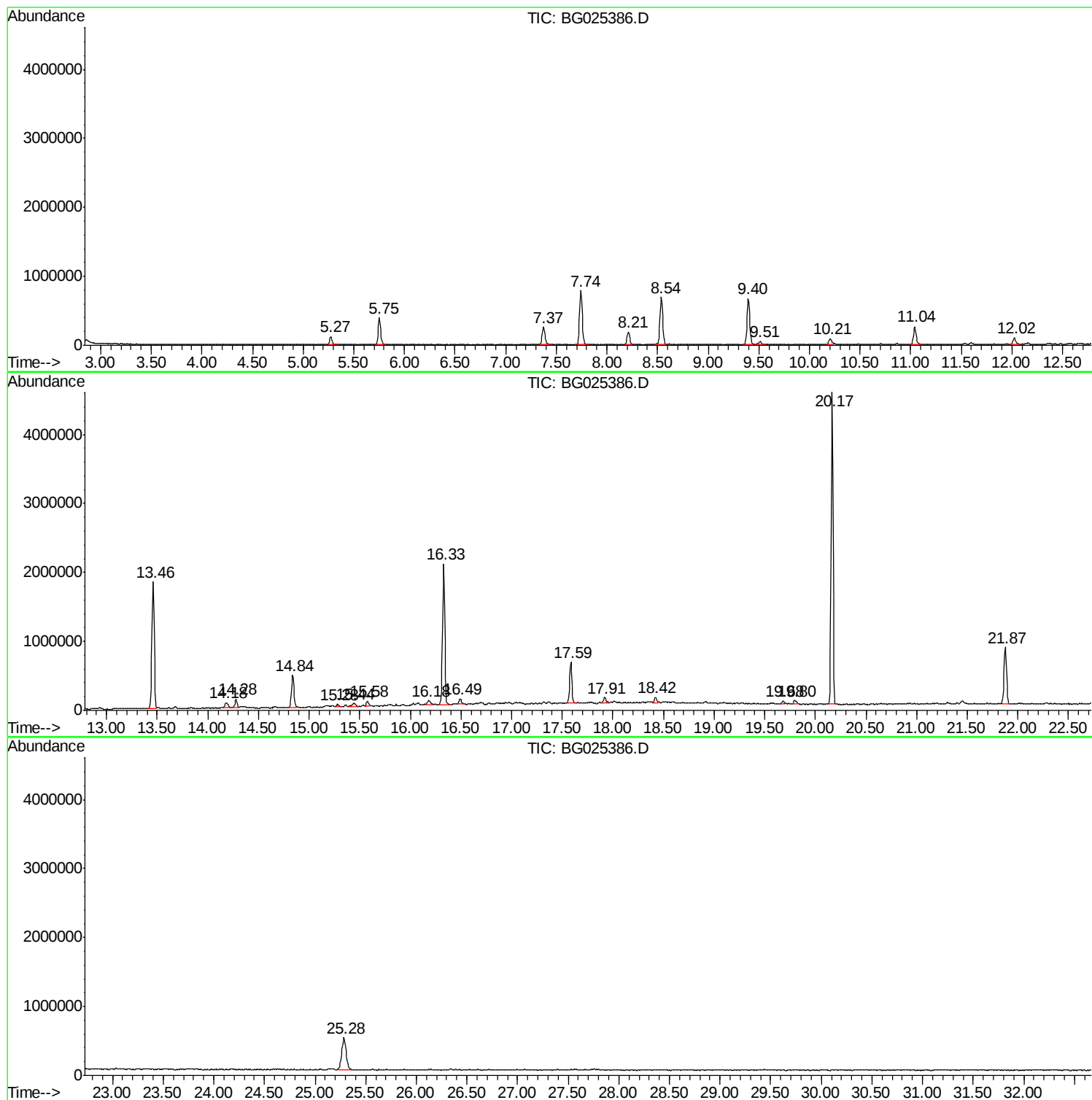
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA G\Methods\8270-BG122116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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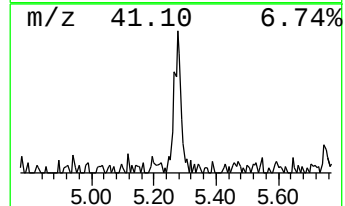
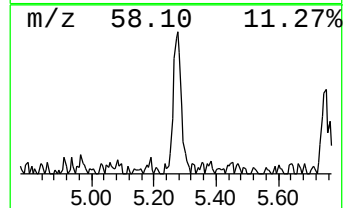
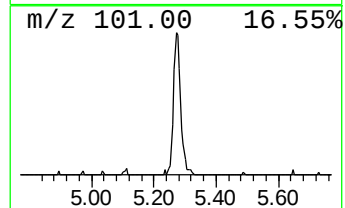
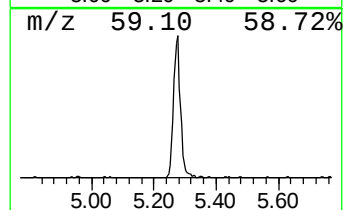
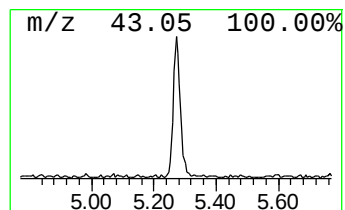
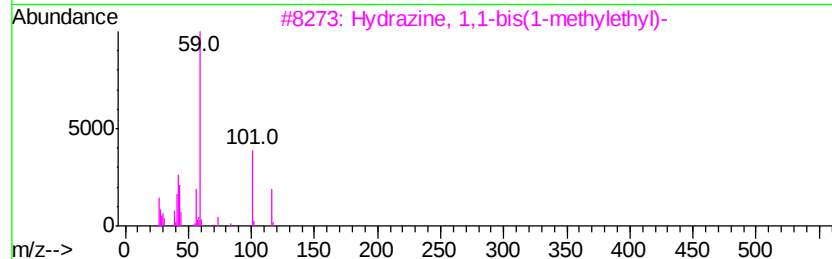
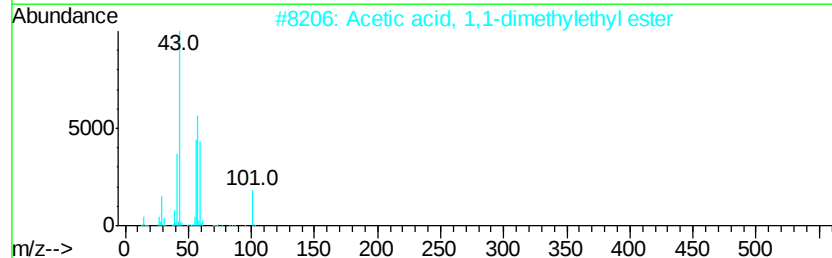
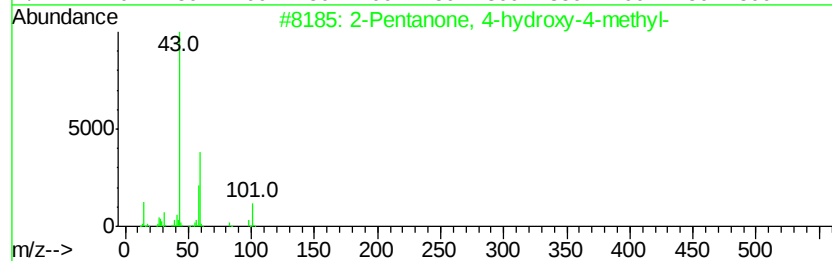
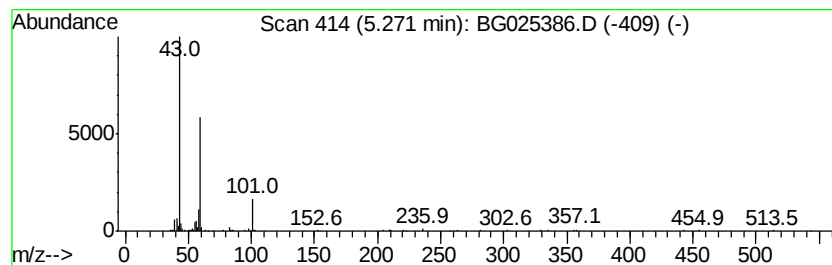
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.27	11.21 ng	193194	1,4-Dichlorobenzene-d4	8.21

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	53
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	38
3			Hydrazine, 1,1-bis(1-methylethyl)-	116	C6H16N2	000921-14-2	37
4			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	27
5			Ethanol, pentamethyl-	116	C7H16O	000594-83-2	25



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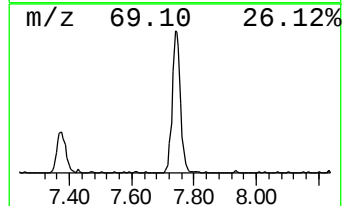
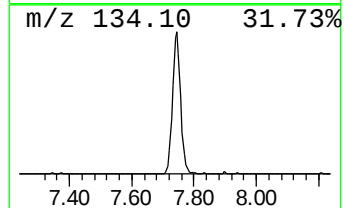
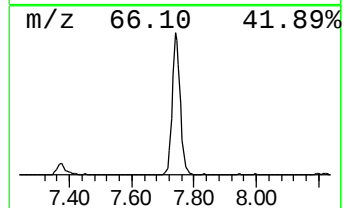
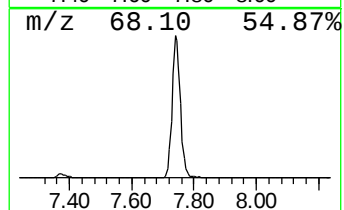
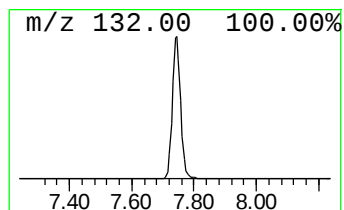
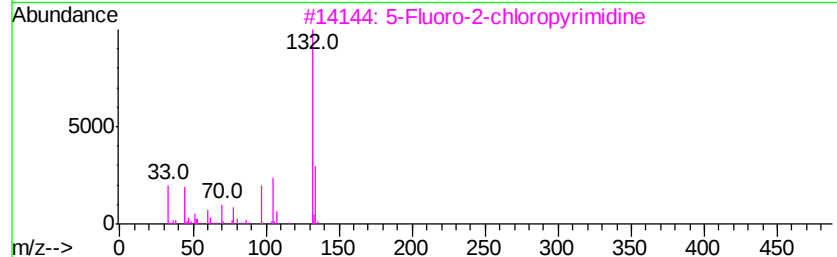
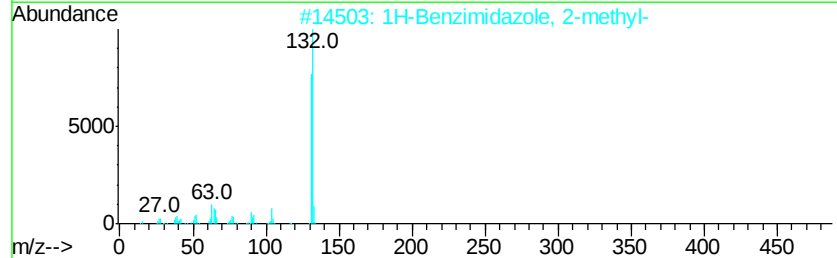
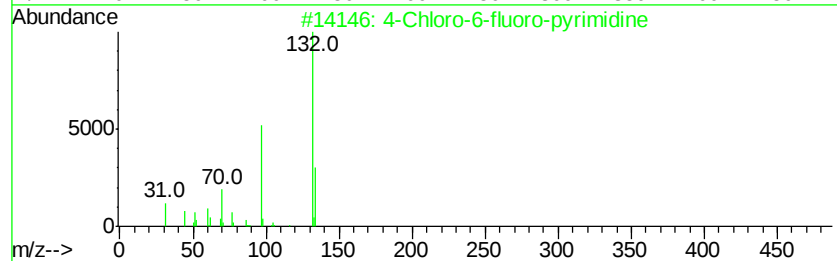
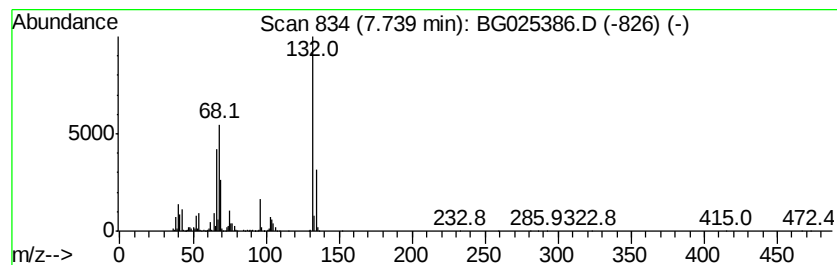
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Peak Number 2 unknown7.74 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.74	84.35 ng	1453680	1,4-Dichlorobenzene-d4	8.21

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	14
2		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	14
3		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	14
4		Cyclopropanamine, 1-phenyl-	133	C9H11N	1000351-26-2	10
5		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	10



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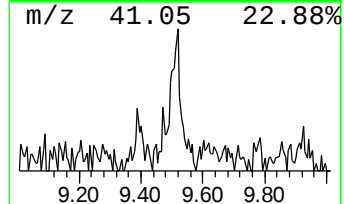
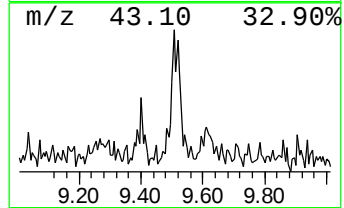
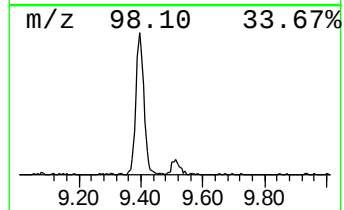
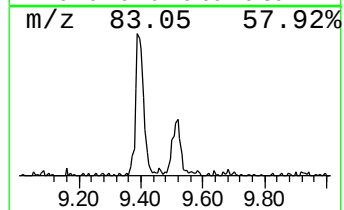
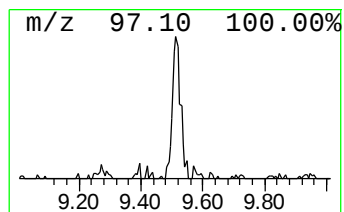
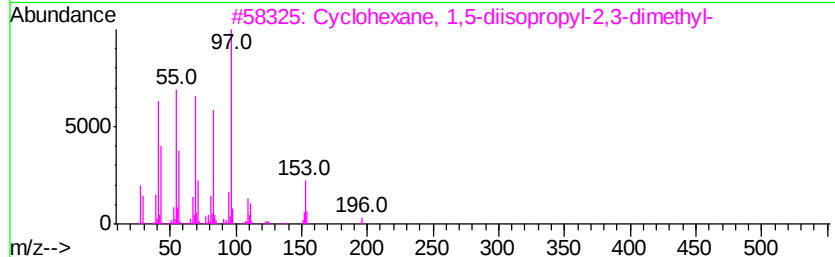
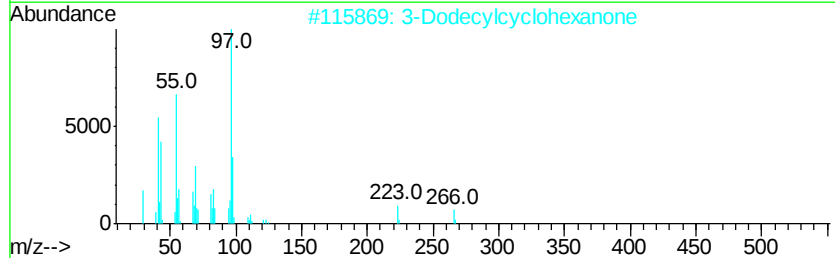
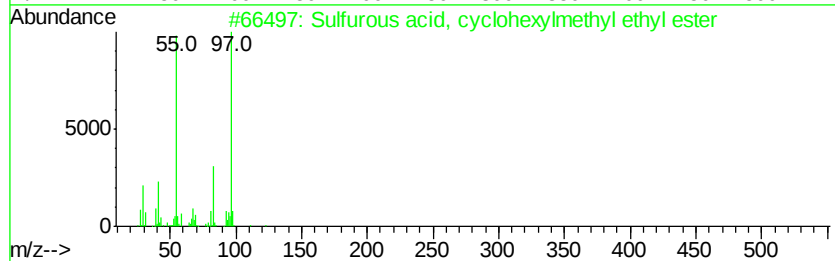
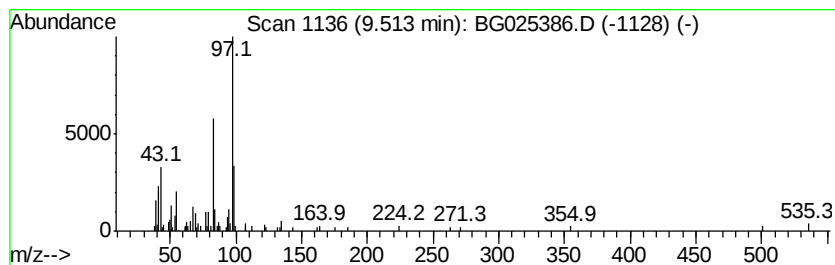
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Peak Number 4 Sulfurous acid, cyclohexylm... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.51	6.50 ng	112022	1,4-Dichlorobenzene-d4	8.21

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Sulfurous acid, cvclohexvlmethylv...	206	C9H18O3S	1000309-21-2	59
2		3-Dodecylcyclohexanone	266	C18H34O	138695-42-8	42
3		Cvclohexane, 1,5-diisopropyl-2,3...	196	C14H28	1000149-58-8	38
4		2-Fluoropvridine	97	C5H4FN	000372-48-5	32
5		Cyclopentanecarboxylic acid, 4-h...	338	C22H42O2	1000282-77-5	28



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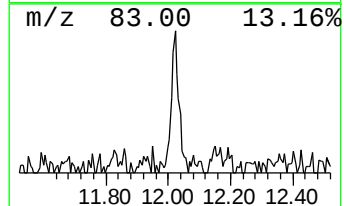
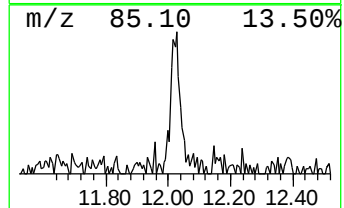
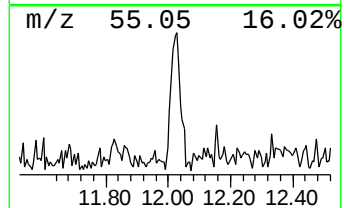
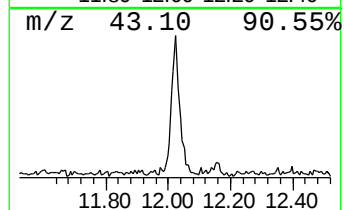
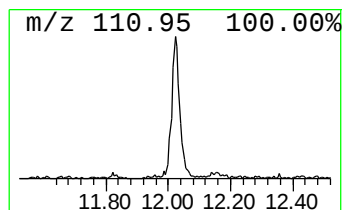
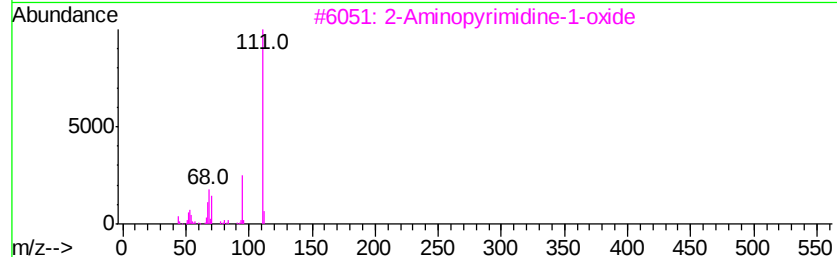
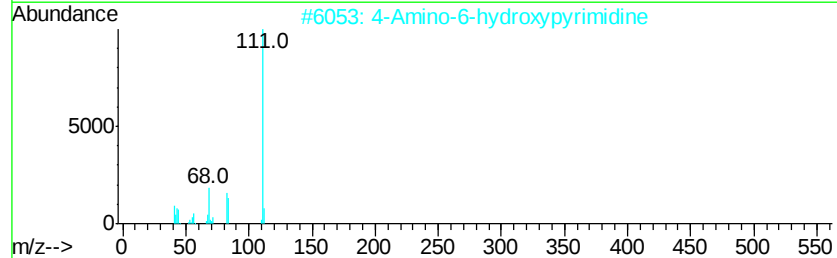
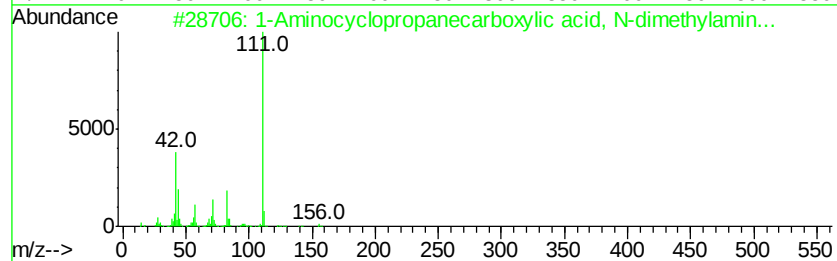
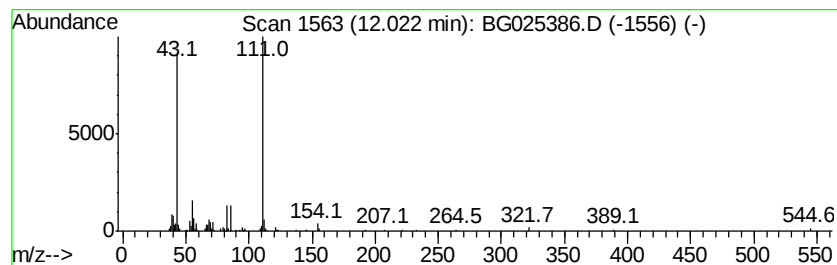
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Peak Number 5 1-Aminocyclopropanecarboxyl... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.02	7.29 ng	184077	Napthalene-d8	11.04

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		1-Aminocyclopropanecarboxylic ac...	156	C7H12N2O2	1000375-50-8	53
2		4-Amino-6-hydroxypyrimidine	111	C4H5N3O	001193-22-2	52
3		2-Aminopyrimidine-1-oxide	111	C4H5N3O	035034-15-2	50
4		2,4-Diamino-1,3,5-triazine	111	C3H5N5	000504-08-5	50
5		Thiophene-2-carboxylic acid, 2-a...	246	C13H10O3S	077708-28-2	50



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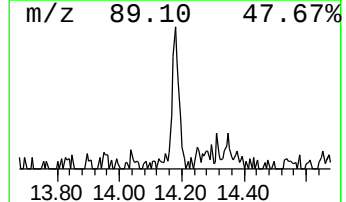
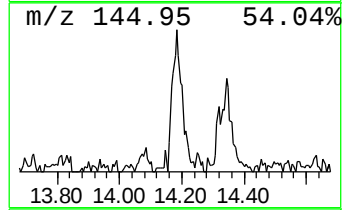
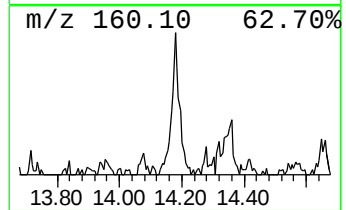
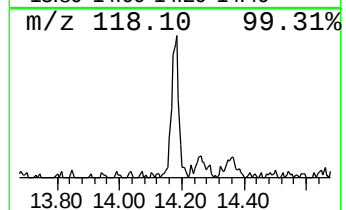
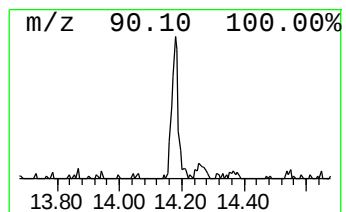
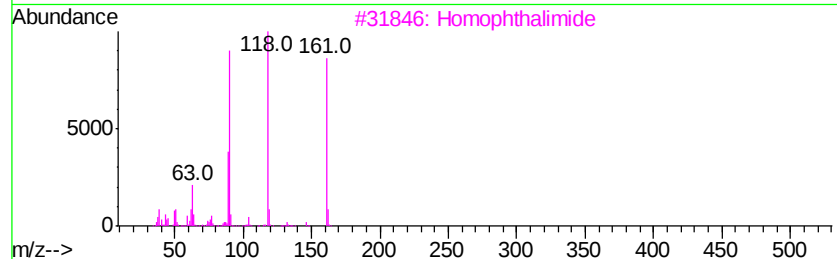
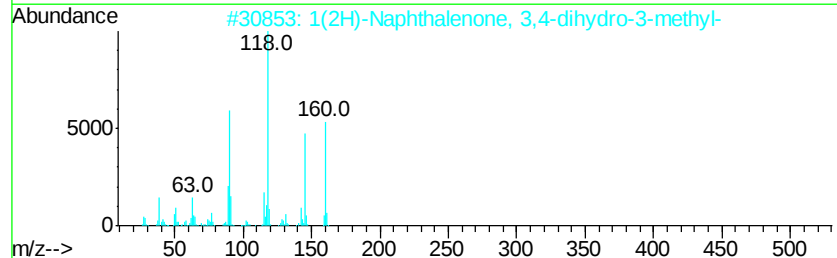
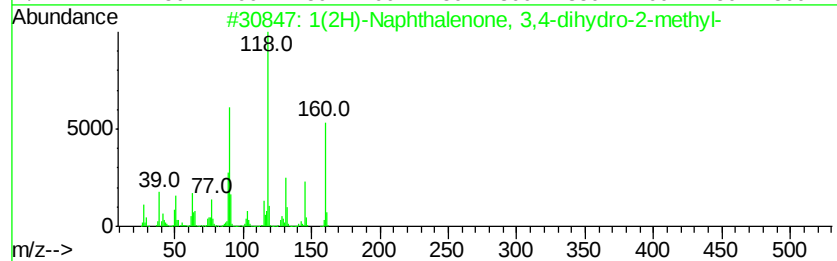
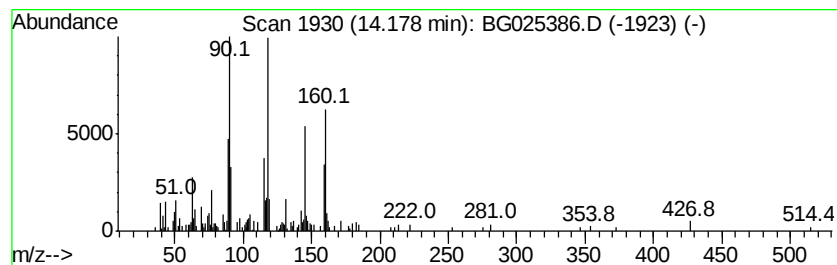
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Peak Number 6 1(2H)-Naphthalenone, 3,4-di... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.18	4.83 ng	194456	Acenaphthene-d10	14.84

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1(2H)-Naphthalenone, 3,4-dihydro...	160	C11H12O	001590-08-5	74
2		1(2H)-Naphthalenone, 3,4-dihydro...	160	C11H12O	014944-23-1	72
3		Homophthalimide	161	C9H7NO2	004456-77-3	58
4		1(2H)-Naphthalenone, 3,4-dihydro-	146	C10H10O	000529-34-0	52
5		4-Methylphthalic anhydride	162	C9H6O3	019438-61-0	50



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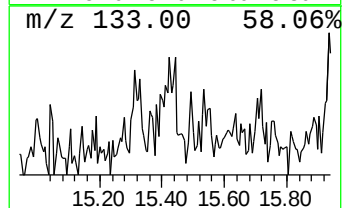
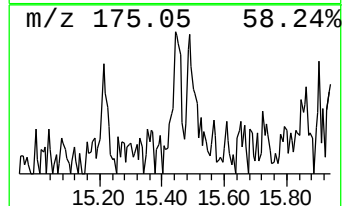
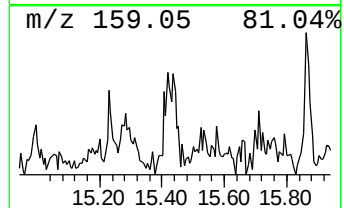
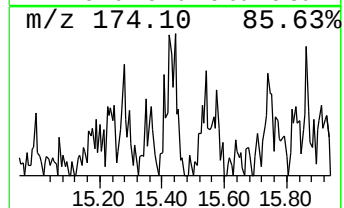
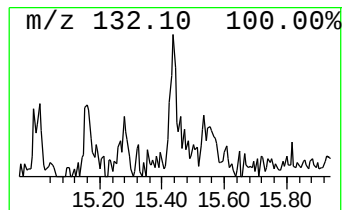
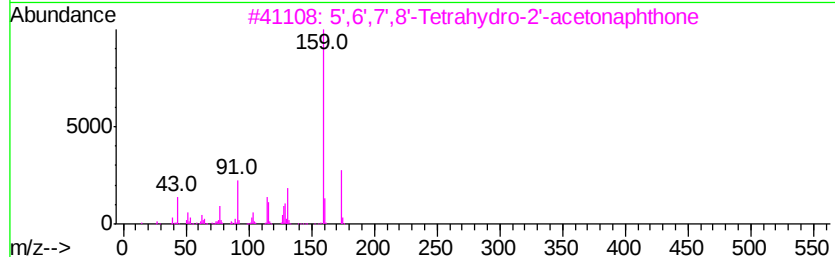
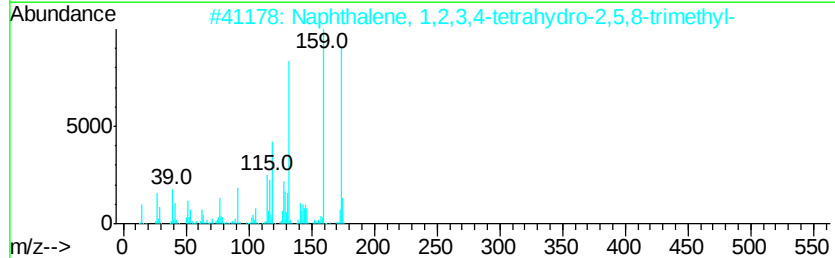
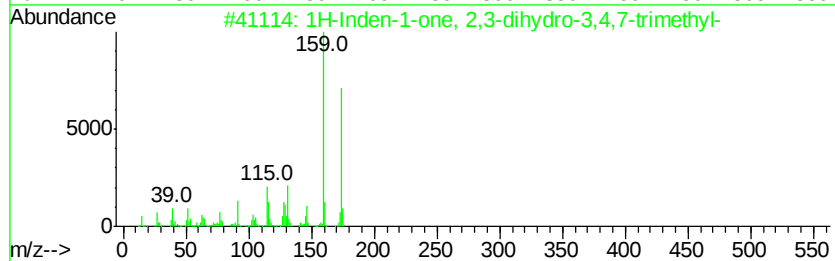
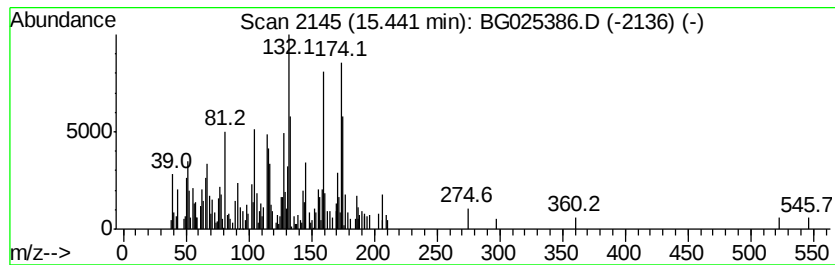
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 1H-Inden-1-one, 2,3-dihydro... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.44	3.37 ng	135612	Acenaphthene-d10	14.84

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		1H-Inden-1-one, 2,3-dihydro-3,4,...	174	C12H14O	035322-84-0	50
2		Naphthalene, 1,2,3,4-tetrahydro-...	174	C13H18	030316-17-7	46
3		5',6',7',8'-Tetrahydro-2'-aceton...	174	C12H14O	000774-55-0	45
4		Tranylcypromine	133	C9H11N	000155-09-9	38
5		Silane, diisothiocyanatodimethyl-	174	C4H6N2S2Si	013125-51-4	38



Data Path : Z:\HPCHEM1\BNA G\DATA\BG122216\
Data File : BG025386.D
Acq On : 22 Dec 2016 12:48
Operator : UM/SJ
Sample : H6199-02
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
W-05

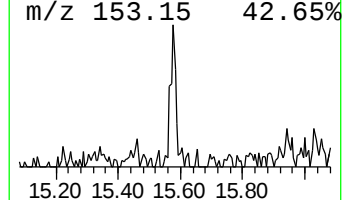
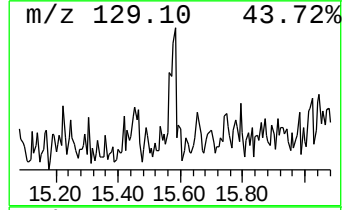
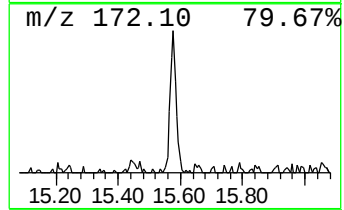
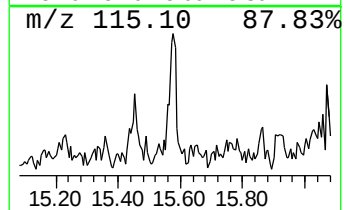
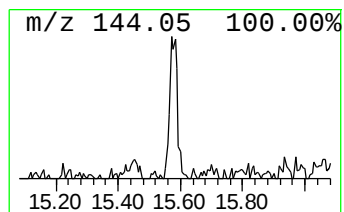
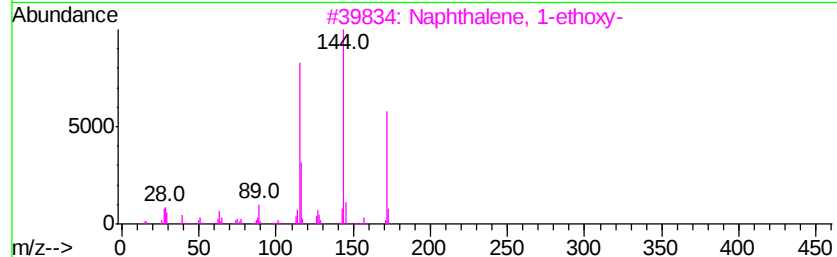
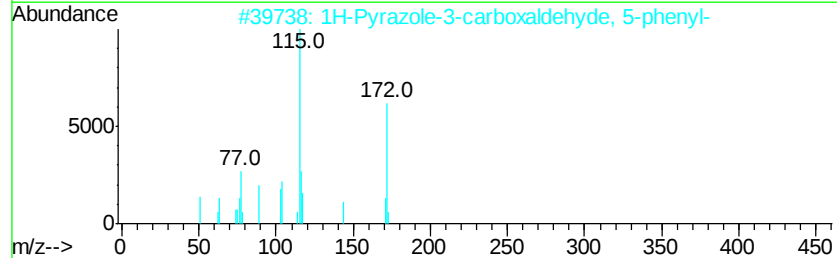
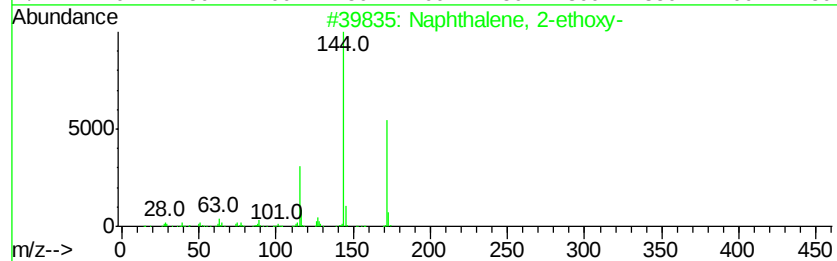
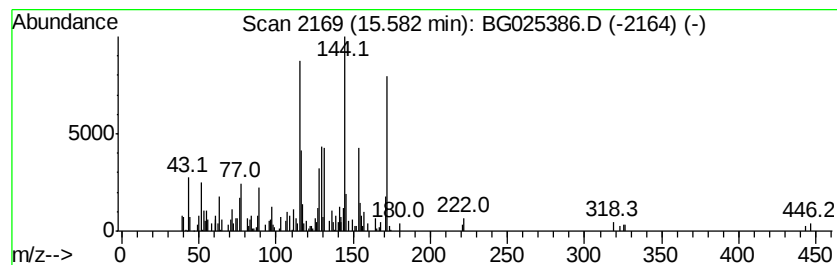
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Naphthalene, 2-ethoxy- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.58	2.80 ng	112683	Acenaphthene-d10	14.84

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 2-ethoxy-	172	C12H12O	000093-18-5	50
2		1H-Pyrazole-3-carboxaldehyde, 5-...	172	C10H8N2O	057204-65-6	46
3		Naphthalene, 1-ethoxy-	172	C12H12O	005328-01-8	45
4		1-Naphthalenecarboxaldehyde, 2-h...	172	C11H8O2	000708-06-5	43
5		2-Naphthalenol	144	C10H8O	000135-19-3	42



Data Path : Z:\HPCHEM1\BNA G\DATA\BG122216\
Data File : BG025386.D
Acq On : 22 Dec 2016 12:48
Operator : UM/SJ
Sample : H6199-02
Misc :
ALS Vial : 21 Sample Multiplier: 1

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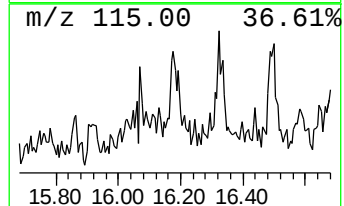
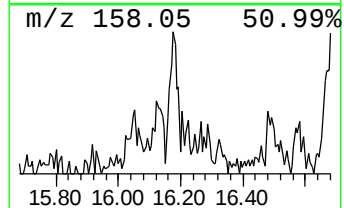
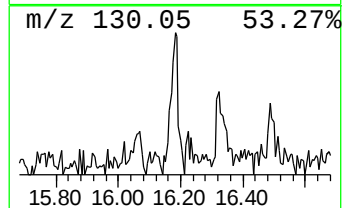
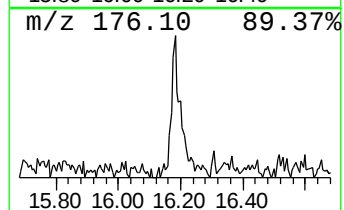
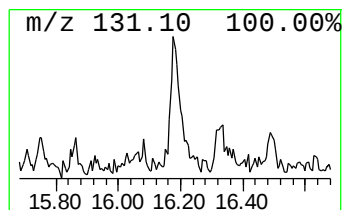
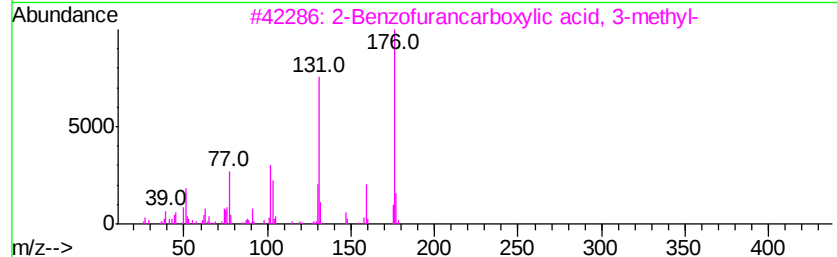
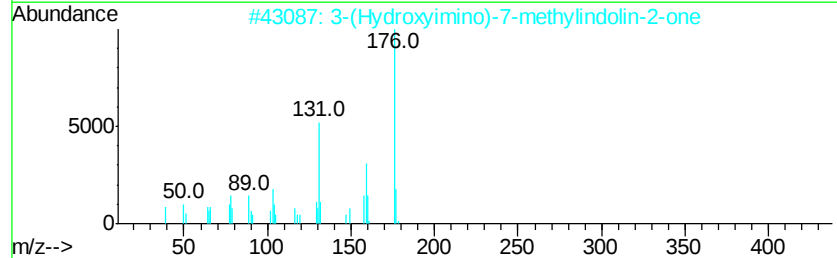
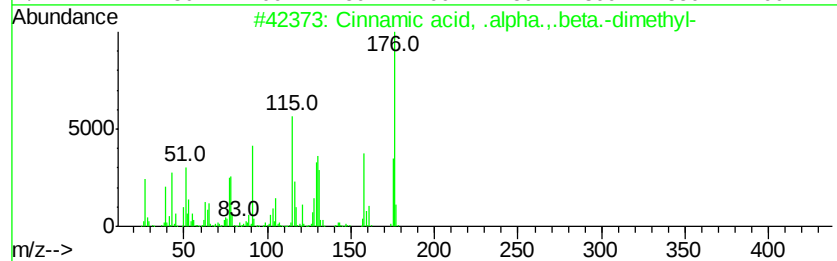
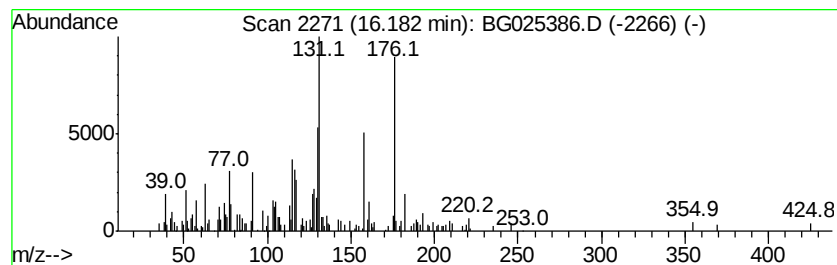
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Cinnamic acid, .alpha.,.bet... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.18	3.07 ng	123651	Acenaphthene-d10	14.84

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cinnamic acid, .alpha.,.beta.-di...	176	C11H12O2	1000151-56-4	53
2		3-(Hydroxyimino)-7-methylindolin...	176	C9H8N2O2	013208-96-3	43
3		2-Benzofurancarboxylic acid, 3-m...	176	C10H8O3	024673-56-1	27
4		1H-Indole, 1-methyl-	131	C9H9N	000603-76-9	27
5		4-Aminoquinaldine	158	C10H10N2	006628-04-2	25



Data Path : Z:\HPCHEM1\BNA G\DATA\BG122216\
Data File : BG025386.D
Acq On : 22 Dec 2016 12:48
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Sample : H6199-02
Misc :
ALS Vial : 21 Sample Multiplier: 1

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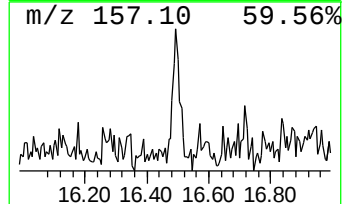
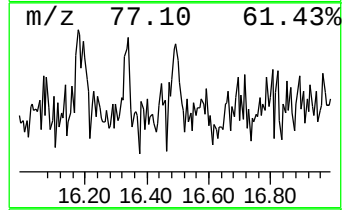
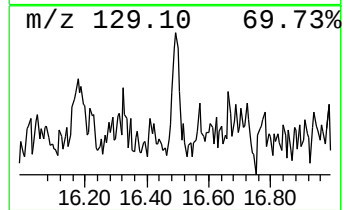
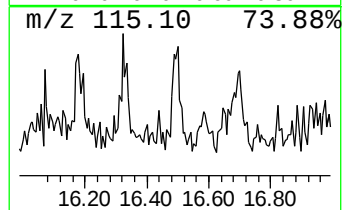
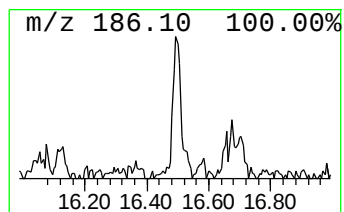
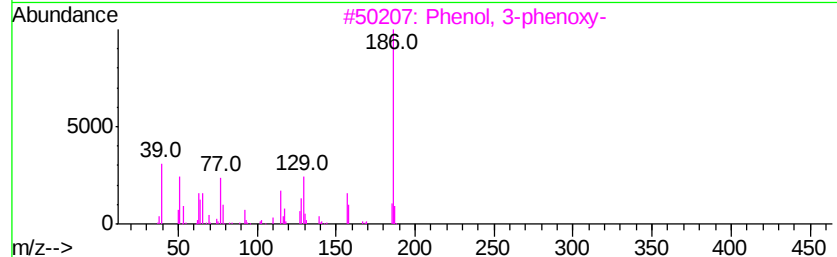
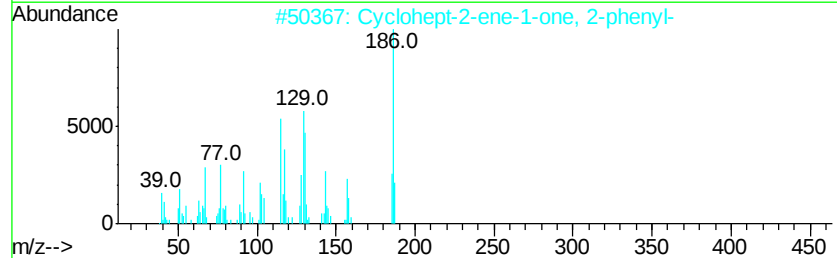
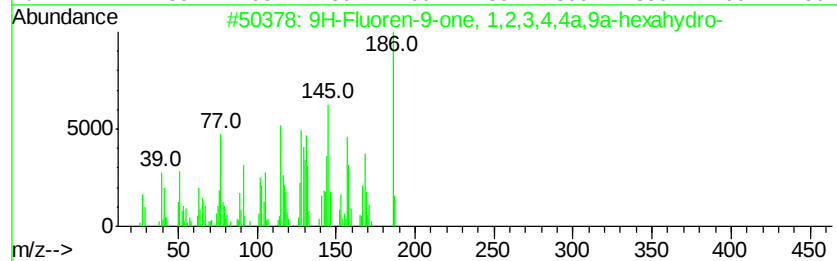
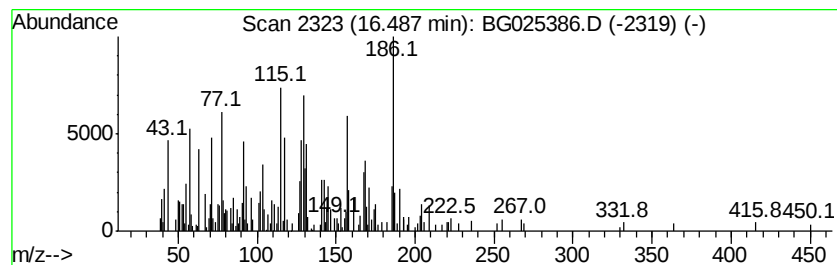
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 9H-Fluoren-9-one, 1,2,3,4,4... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.49	3.07 ng	143581	Phenanthrene-d10	17.59

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	9H-Fluoren-9-one, 1,2,3,4,4a,9a-...	186	C13H14O	001203-67-4	68
2		Cyclohept-2-ene-1-one, 2-phenyl-	186	C13H14O	065825-75-4	55
3		Phenol, 3-phenoxy-	186	C12H10O2	000713-68-8	50
4		5-Thiazolecarboxylic acid, 2-amino-	186	C7H10N2O2S	007210-76-6	42
5		[1,1'-Biphenyl]-4,4'-diol	186	C12H10O2	000092-88-6	42



Data Path : Z:\HPCHEM1\BNA G\DATA\BG122216\
Data File : BG025386.D
Acq On : 22 Dec 2016 12:48
Operator : UM/SJ
Sample : H6199-02
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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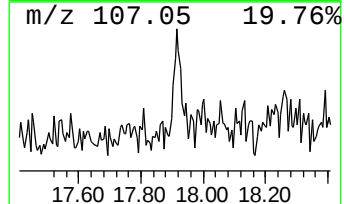
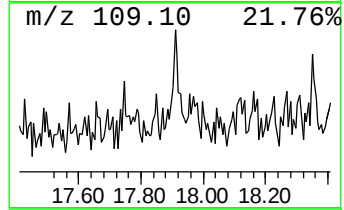
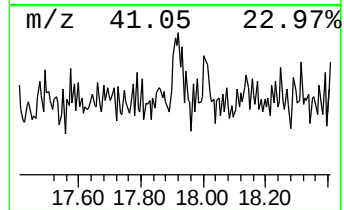
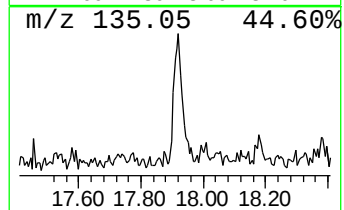
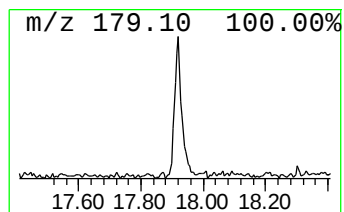
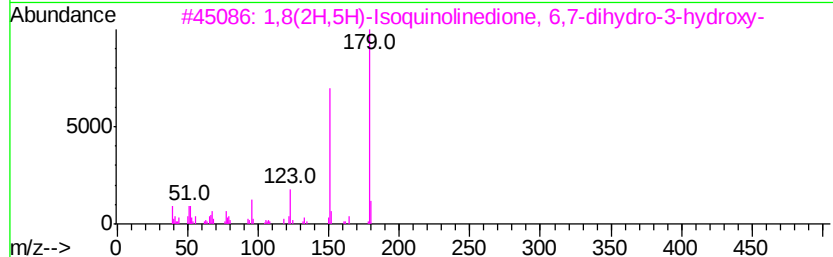
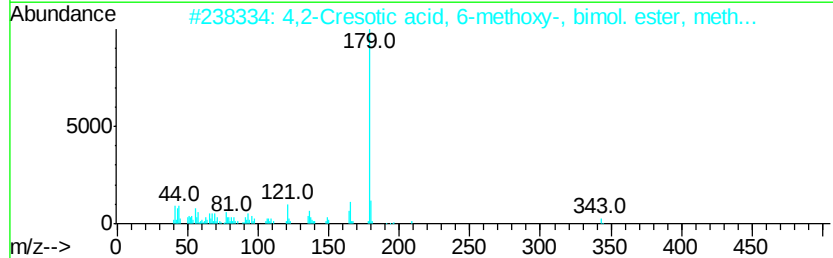
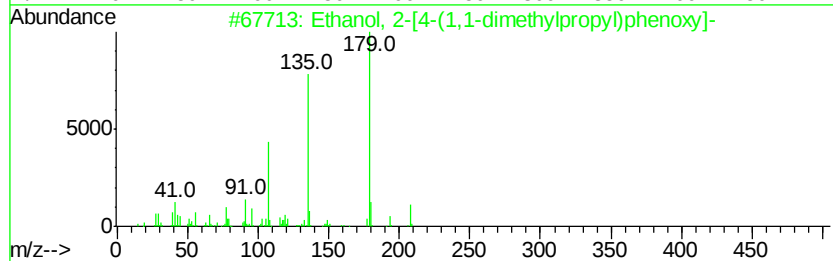
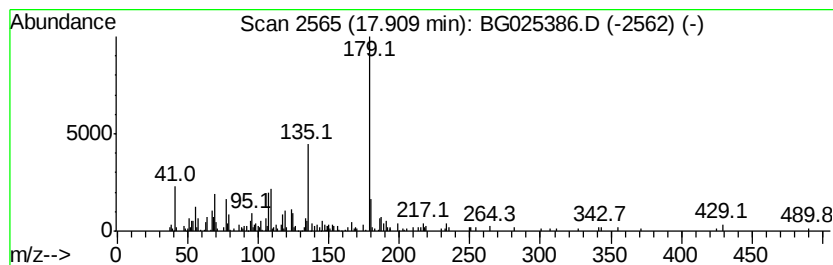
Quant Method : Z:\HPCHEM1\BNA G\Methods\8270-BG122116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Ethanol, 2-[4-(1,1-dimethyl... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.91	2.77 ng	129293	Phenanthrene-d10	17.59

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Ethanol, 2-[4-(1,1-dimethylpropyl)phenoxy]-	208	C13H20O2	006382-07-6	64
2		4,2-Cresotic acid, 6-methoxy-, bimol. ester, meth...	538	C29H30O10	019314-74-0	47
3		1,8(2H,5H)-Isoquinolinedione, 6,7-dihydro-3-hydroxy-	179	C9H9NO3	037704-54-4	47
4		4-Butylbenzoic acid, tetradecyl ester, meth...	374	C25H42O2	1000292-32-7	43
5		2,4-Dimethoxyphenyl isocyanate	179	C9H9NO3	084370-87-6	43



Data Path : Z:\HPCHEM1\BNA G\DATA\BG122216\
Data File : BG025386.D
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Sample : H6199-02
Misc :
ALS Vial : 21 Sample Multiplier: 1

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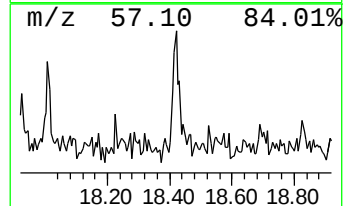
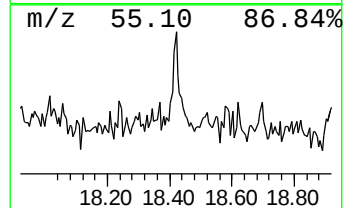
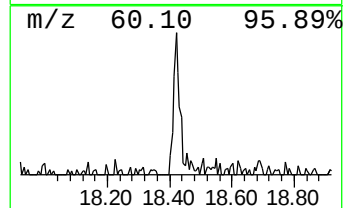
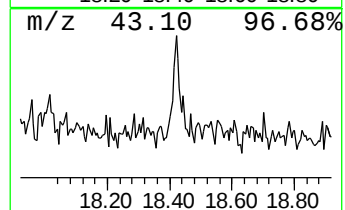
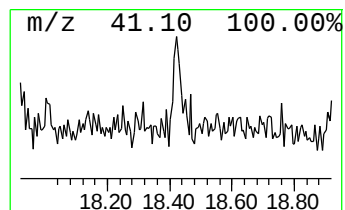
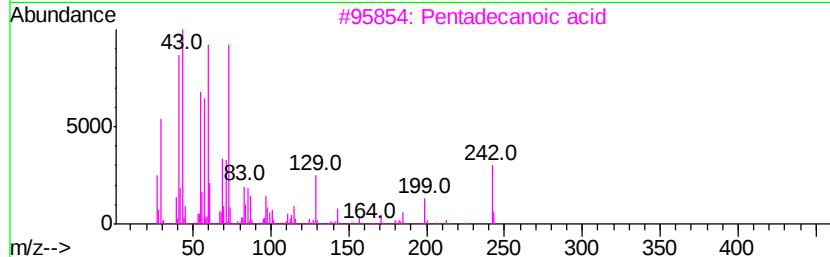
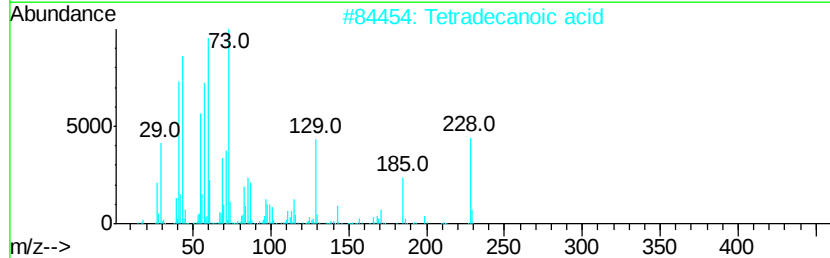
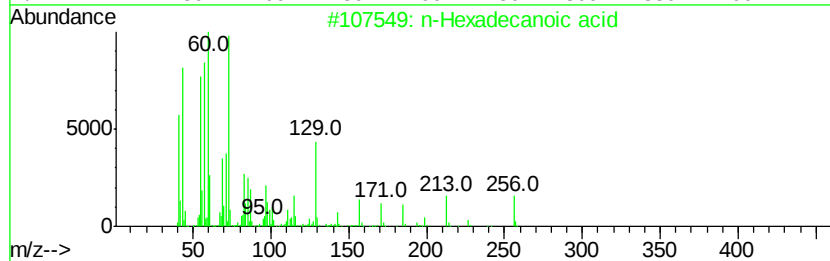
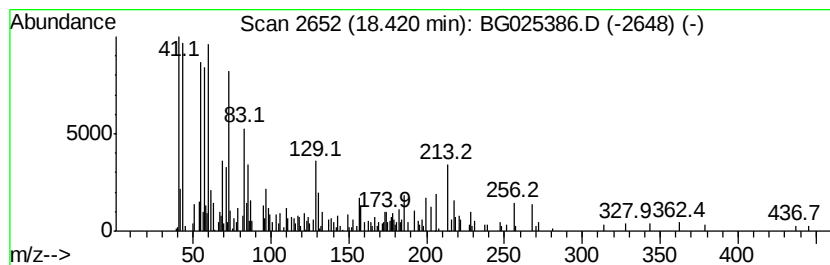
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 n-Hexadecanoic acid Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.42	2.58 ng	120399	Phenanthrene-d10	17.59

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	91
2			Tetradecanoic acid	228	C14H28O2	000544-63-8	87
3			Pentadecanoic acid	242	C15H30O2	001002-84-2	43
4			Tridecanoic acid	214	C13H26O2	000638-53-9	43
5			Dodecanoic acid	200	C12H24O2	000143-07-7	38



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG122216\
Data File : BG025386.D
Acq On : 22 Dec 2016 12:48
Operator : UM/SJ
Sample : H6199-02
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_G\Methods\8270-BG122116.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
2-Pentanone, 4-hy...	5.27	11.2	ng	193194	1	8.21	344696	20.0
unknown7.74	7.74	84.3	ng	1453680	1	8.21	344696	20.0
Sulfurous acid, c...	9.51	6.5	ng	112022	1	8.21	344696	20.0
1-Aminocyclopropa...	12.02	7.3	ng	184077	2	11.04	505099	20.0
1(2H)-Naphthaleno...	14.18	4.8	ng	194456	3	14.84	804666	20.0
1H-Inden-1-one, 2...	15.44	3.4	ng	135612	3	14.84	804666	20.0
Naphthalene, 2-et...	15.58	2.8	ng	112683	3	14.84	804666	20.0
Cinnamic acid, .a...	16.18	3.1	ng	123651	3	14.84	804666	20.0
9H-Fluoren-9-one, ...	16.49	3.1	ng	143581	4	17.59	934989	20.0
Ethanol, 2-[4-(1,...	17.91	2.8	ng	129293	4	17.59	934989	20.0
n-Hexadecanoic acid	18.42	2.6	ng	120399	4	17.59	934989	20.0