

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG041715\
 Data File : BG016545.D
 Acq On : 19 Apr 2015 5:30
 Operator : TP/IZ
 Sample : G1765-02
 Misc :
 ALS Vial : 54 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 AS-2

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-BG040615.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	2.748	5	10	20	rVB2	94050	178571	9.26%	0.993%
2	2.831	20	24	29	rBV	126671	169835	8.81%	0.944%
3	2.925	35	40	52	rVB	189492	321335	16.66%	1.786%
4	4.770	346	354	364	rBV	280284	411065	21.31%	2.285%
5	5.251	429	436	447	rBV	782122	1141526	59.18%	6.346%
6	6.844	694	707	720	rBV	763920	1243181	64.45%	6.911%
7	7.196	760	767	777	rVB	833296	1312878	68.07%	7.298%
8	7.654	836	845	852	rBV	330001	522019	27.06%	2.902%
9	7.978	891	900	908	rVB	670909	1062095	55.06%	5.904%
10	8.812	1036	1042	1050	rBV	479529	780634	40.47%	4.340%
11	10.439	1307	1319	1328	rBV	473435	840353	43.57%	4.672%
12	12.919	1731	1741	1751	rBV	1285565	1928810	100.00%	10.722%
13	13.712	1872	1876	1881	rBV	204790	272931	14.15%	1.517%
14	13.830	1892	1896	1905	rVB5	67615	139959	7.26%	0.778%
15	13.953	1913	1917	1921	rBV3	105108	132823	6.89%	0.738%
16	14.035	1926	1931	1936	rBV4	179816	332458	17.24%	1.848%
17	14.123	1942	1946	1954	rVB	310960	471845	24.46%	2.623%
18	14.200	1954	1959	1964	rBV4	109888	198608	10.30%	1.104%
19	14.264	1965	1970	1972	rBV	166903	282404	14.64%	1.570%
20	14.300	1972	1976	1985	rVV2	702010	1189913	61.69%	6.615%
21	14.388	1987	1991	1996	rVB4	71670	129897	6.73%	0.722%
22	14.828	2061	2066	2073	rBV7	104454	247193	12.82%	1.374%
23	15.016	2094	2098	2105	rBV7	115264	266153	13.80%	1.480%
24	15.087	2105	2110	2114	rBV2	379608	582552	30.20%	3.238%
25	15.792	2226	2230	2236	rVB2	573823	843156	43.71%	4.687%
26	16.039	2265	2272	2279	rBV2	420203	1015586	52.65%	5.646%
27	16.861	2409	2412	2422	rVB	502867	843206	43.72%	4.687%
28	17.043	2440	2443	2448	rBV2	769466	1127498	58.46%	6.268%

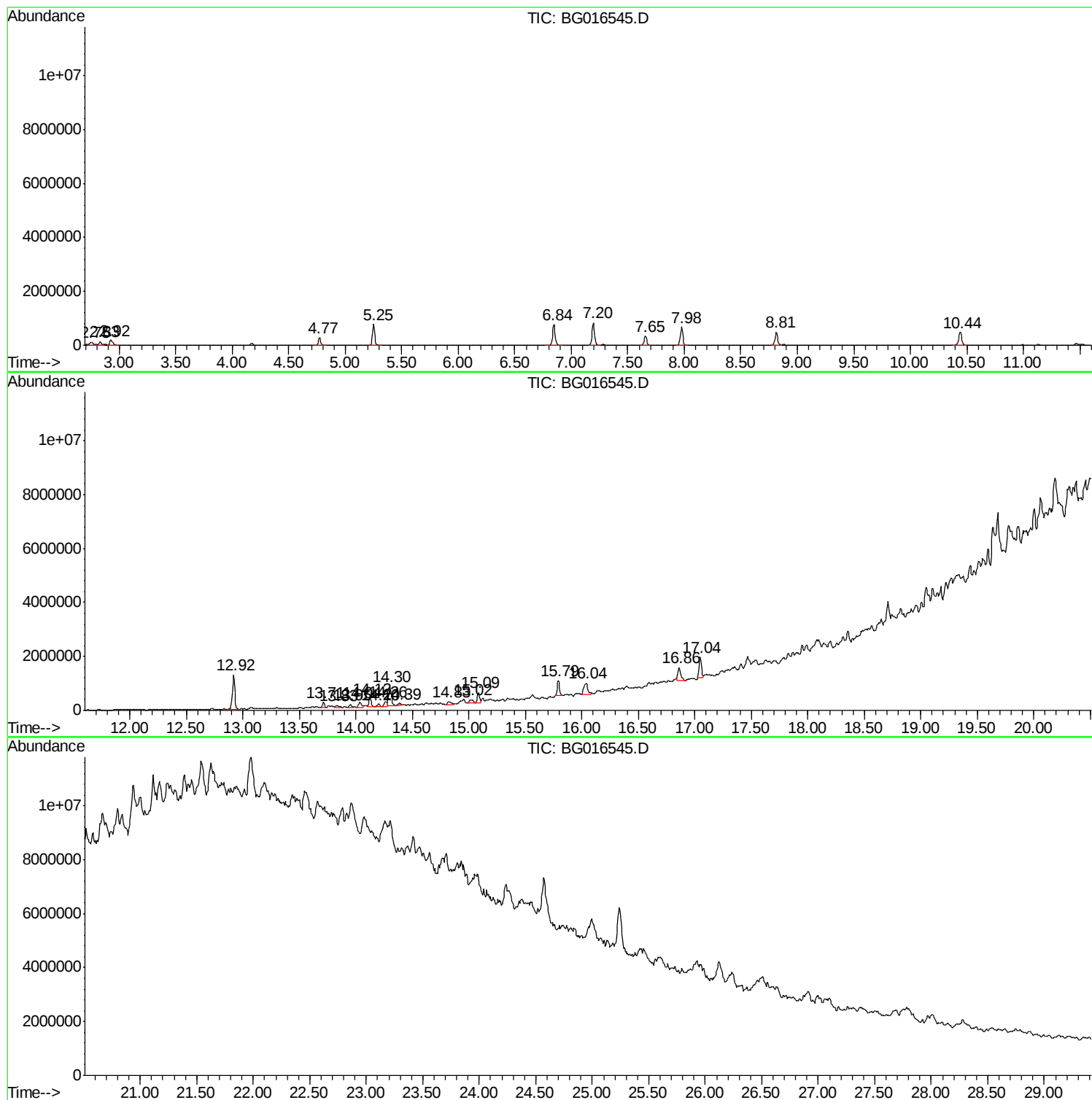
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Instrument :
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Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG040615.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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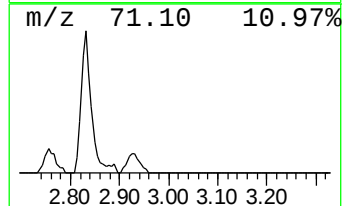
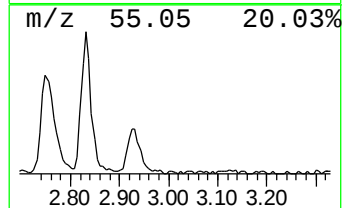
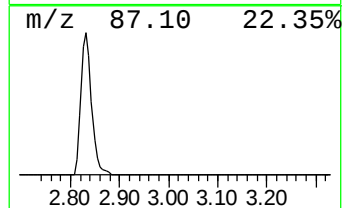
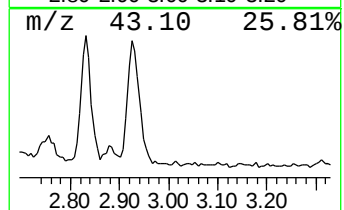
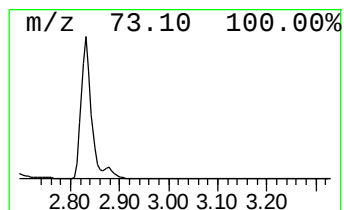
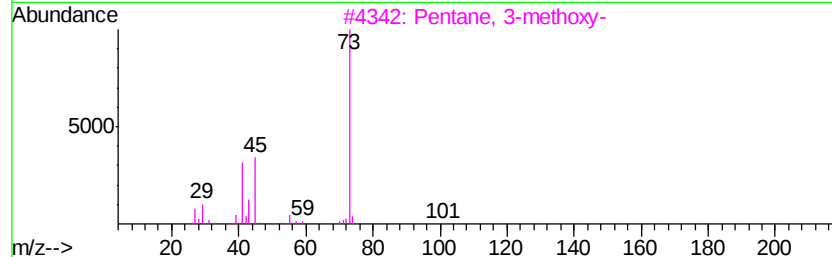
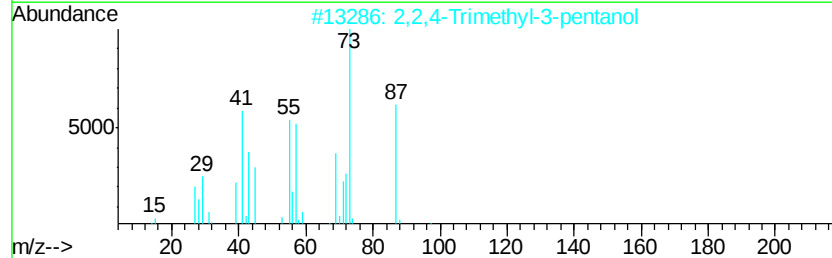
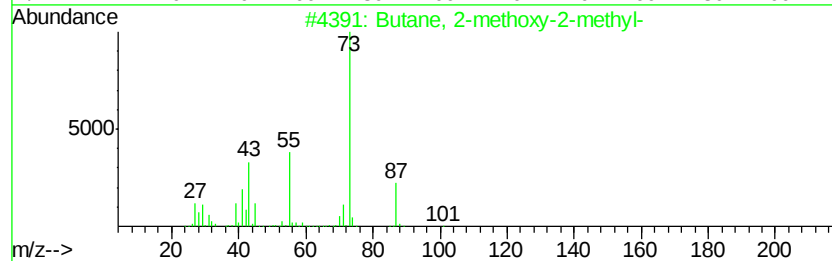
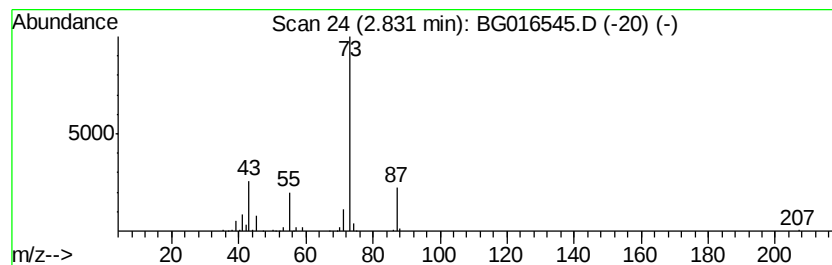
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 Butane, 2-methoxy-2-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.83	6.51 ng	169835	1,4-Dichlorobenzene-d4	7.65

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	64
3		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	17
4		1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	9
5		N-Ethylformamide	73	C3H7NO	000627-45-2	9



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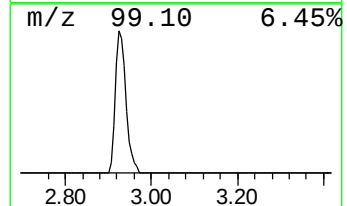
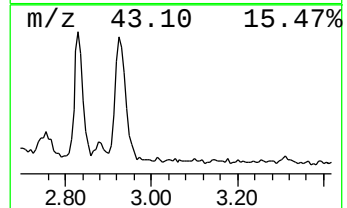
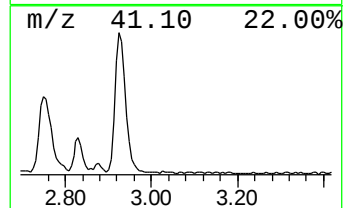
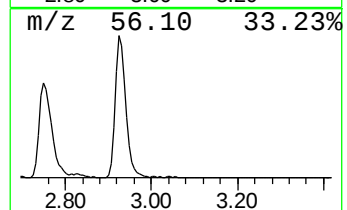
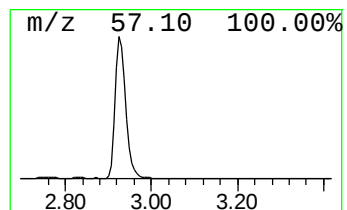
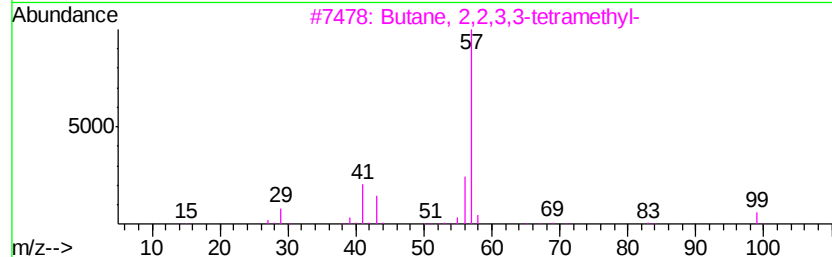
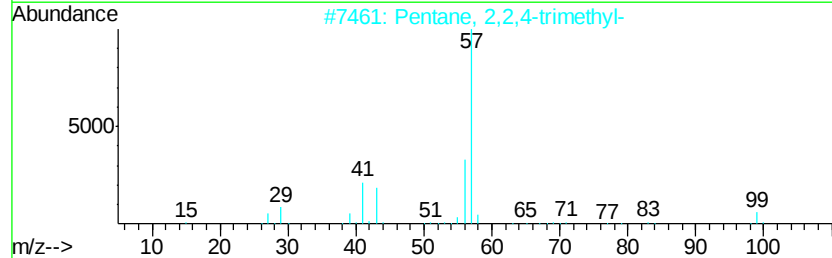
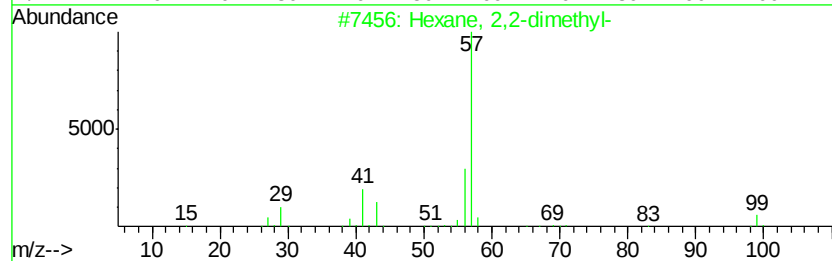
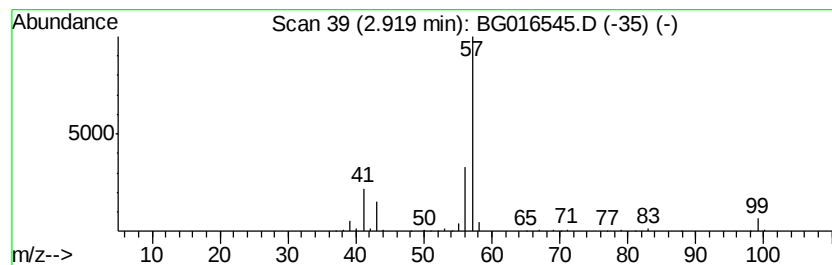
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Hexane, 2,2-dimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.92	12.31 ng	321335	1,4-Dichlorobenzene-d4	7.65

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexane, 2,2-dimethyl-	114	C8H18	000590-73-8	83
2		Pentane, 2,2,4-trimethyl-	114	C8H18	000540-84-1	83
3		Butane, 2,2,3,3-tetramethyl-	114	C8H18	000594-82-1	83
4		Pentane, 2,2,3-trimethyl-	114	C8H18	000564-02-3	64
5		Heptane, 2,2,4,6,6-pentamethyl-	170	C12H26	013475-82-6	64



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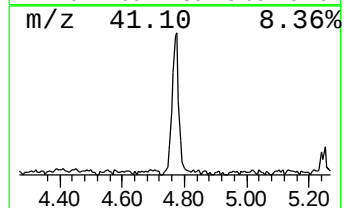
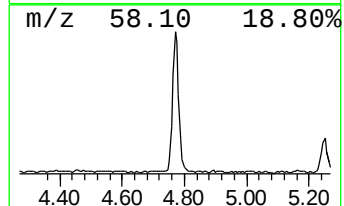
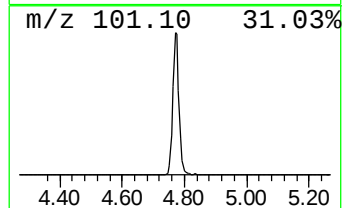
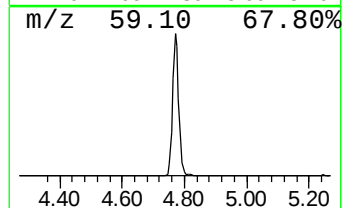
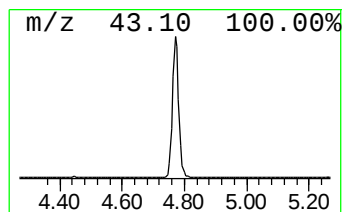
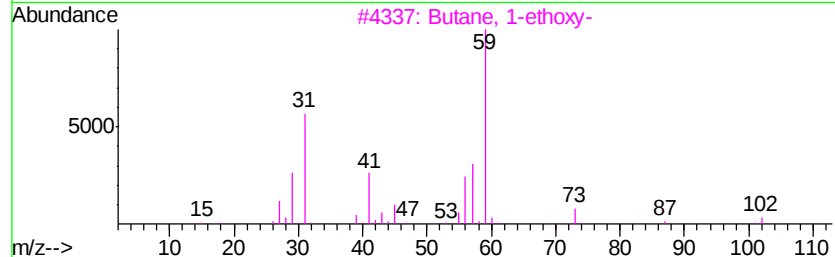
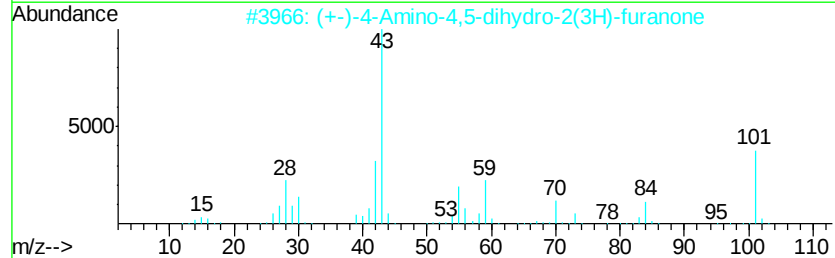
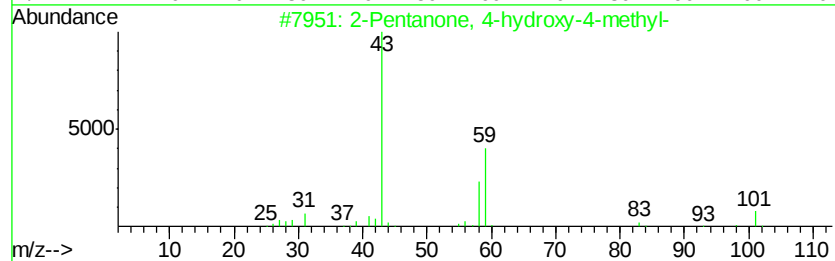
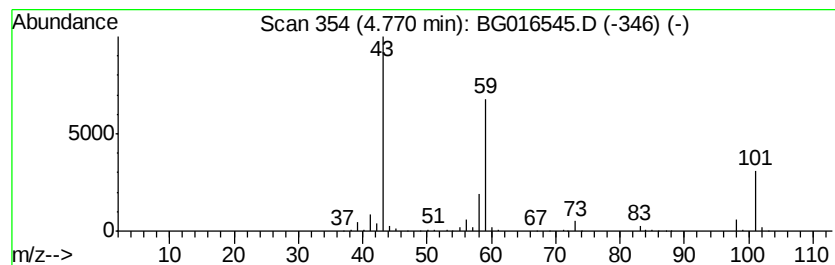
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.77	15.75 ng	411065	1,4-Dichlorobenzene-d4	7.65

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2		(+)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	47
3		Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9
4		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
5		5-Hexen-2-one	98	C6H10O	000109-49-9	9



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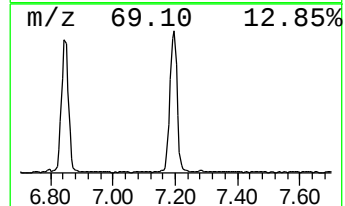
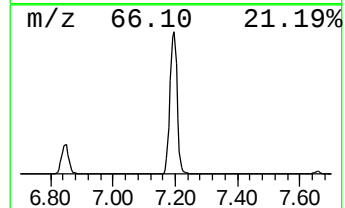
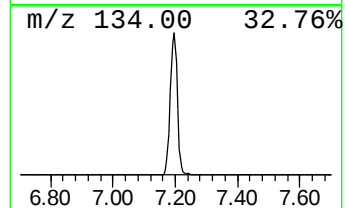
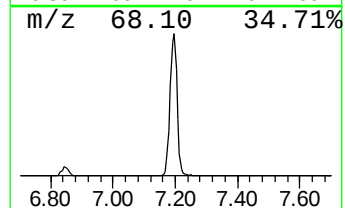
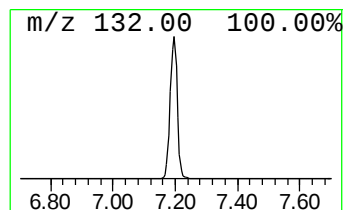
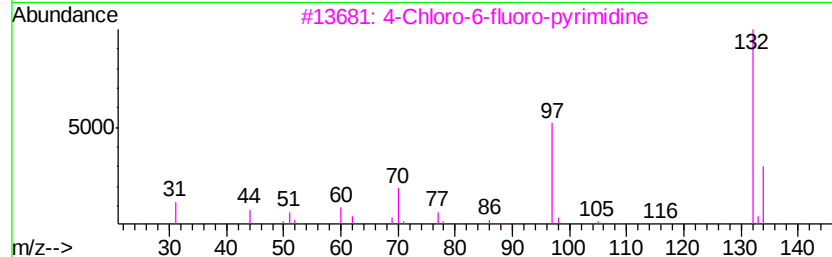
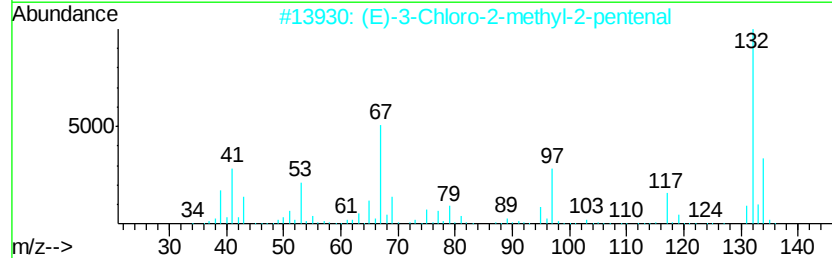
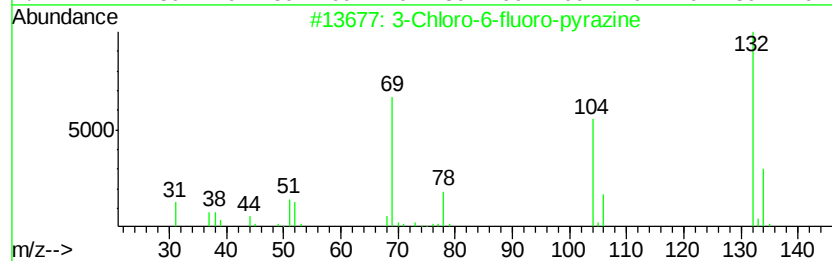
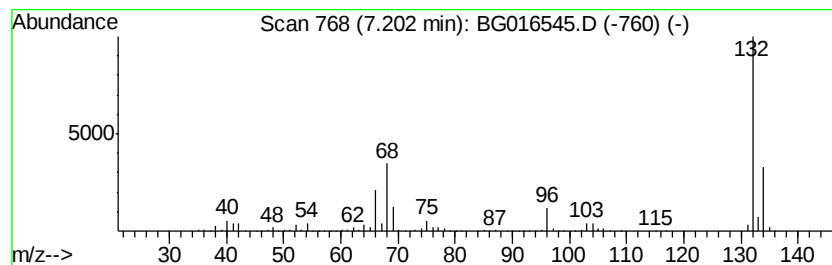
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Peak Number 5 unknown7.20 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.20	50.30 ng	1312880	1,4-Dichlorobenzene-d4	7.65

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	38
2	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	35
3	4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	9
4	2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9
5	4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9



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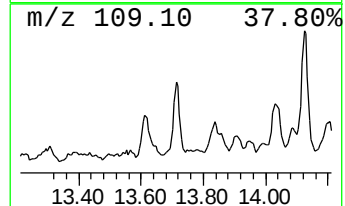
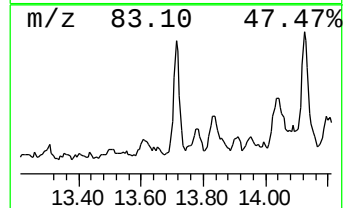
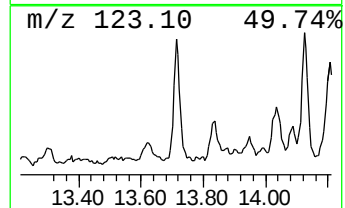
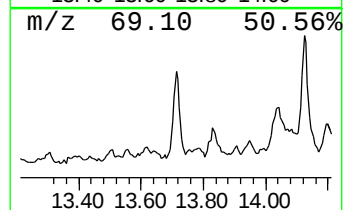
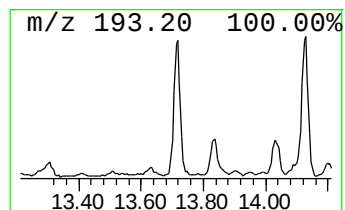
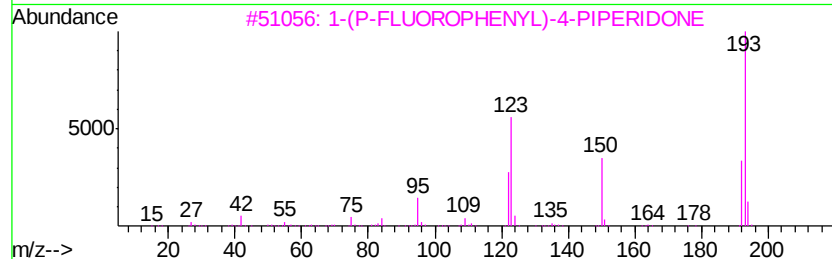
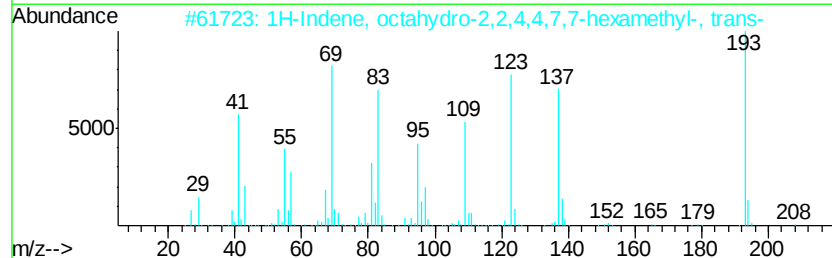
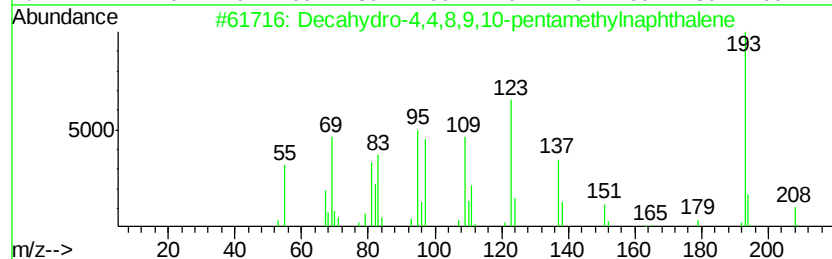
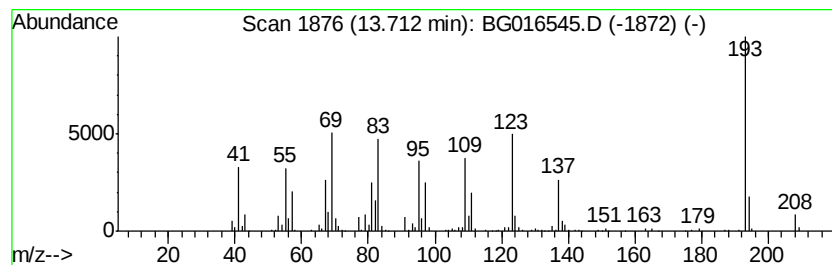
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 7 Decahydro-4,4,8,9,10-pentam... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.71	4.59 ng	272931	Acenaphthene-d10	14.30

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Decahydro-4,4,8,9,10-pentamethyl...	208	C15H28	080655-44-3	83
2		1H-Indene, octahydro-2,2,4,4,7,7...	208	C15H28	054832-83-6	40
3		1-(P-FLUOROPHENYL)-4-PIPERIDONE	193	C11H12FNO	1000238-56-7	37
4		2-Anthracenamine	193	C14H11N	000613-13-8	30
5		Iminostilbene	193	C14H11N	000256-96-2	22



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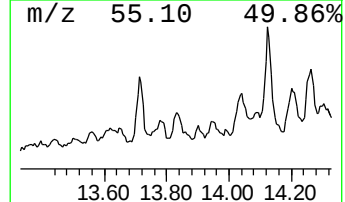
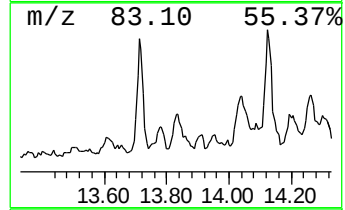
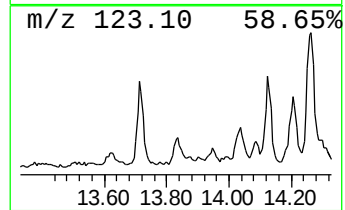
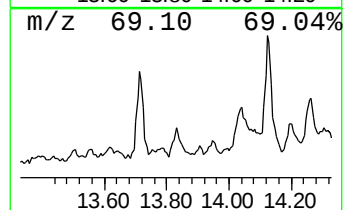
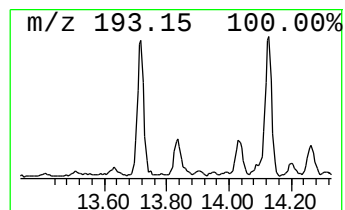
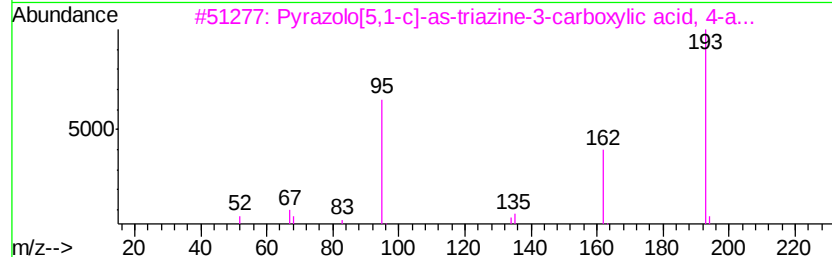
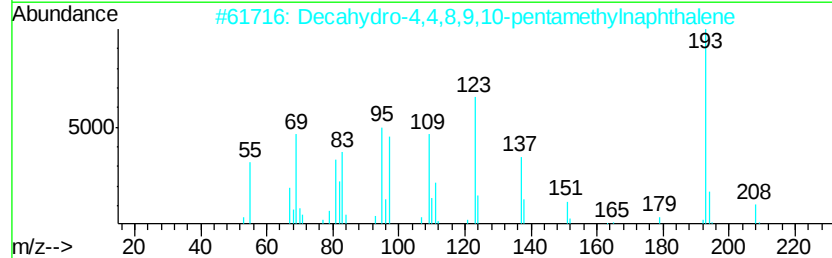
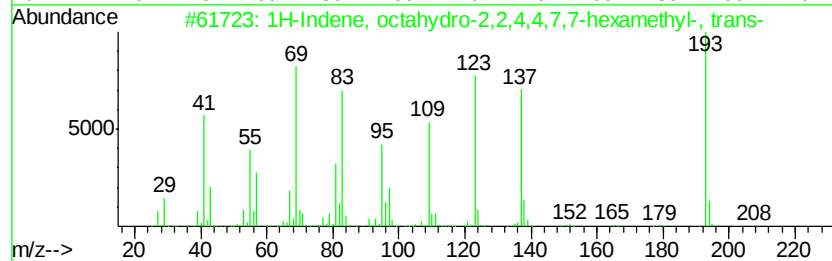
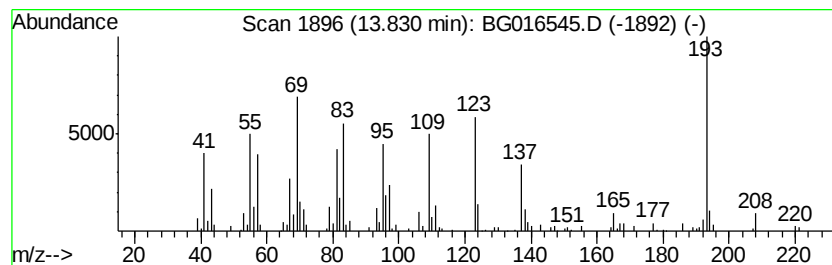
Instrument :
BNA_G
ClientSampleId :
AS-2

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG040615.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 1H-Indene, octahydro-2,2,4,... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.83	2.35 ng	139959	Acenaphthene-d10	14.30
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	1H-Indene, octahydro-2,2,4,4,7,7...	208 C15H28	054832-83-6	80
2	Decahydro-4,4,8,9,10-pentamethyl...	208 C15H28	080655-44-3	45
3	Pyrazolo[5,1-c]-as-triazine-3-ca...	193 C6H7N7O	016111-78-7	35
4	Acridine, 9-methyl-	193 C14H11N	000611-64-3	18
5	Borinic acid, diethyl-, 1-cyclod...	250 C16H31BO	061142-73-2	16



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG041715\
Data File : BG016545.D
Acq On : 19 Apr 2015 5:30
Operator : TP/IZ
Sample : G1765-02
Misc :
ALS Vial : 54 Sample Multiplier: 1

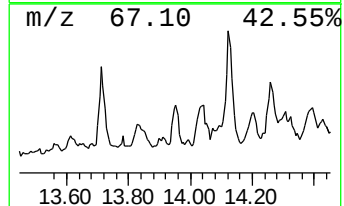
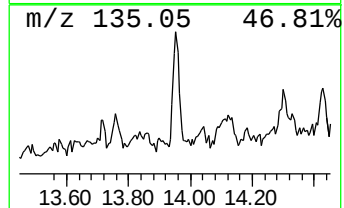
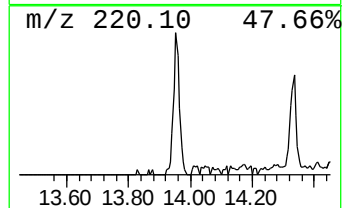
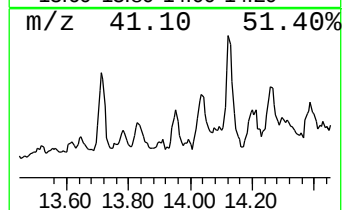
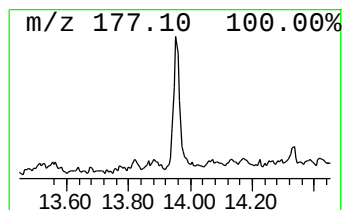
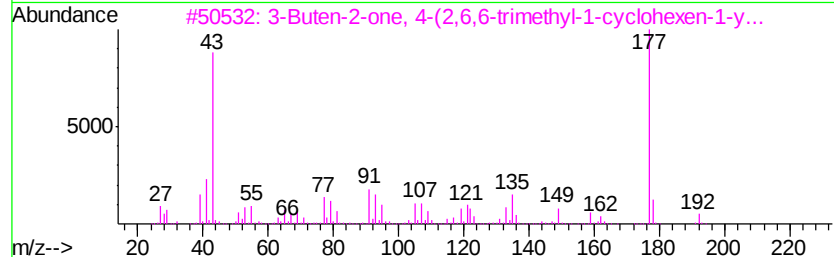
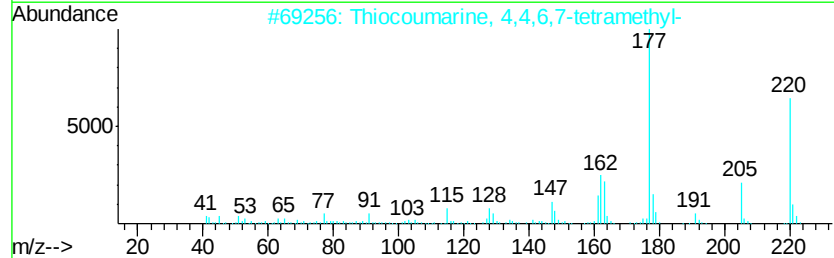
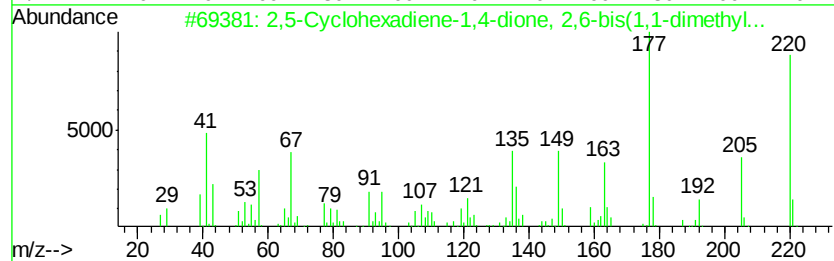
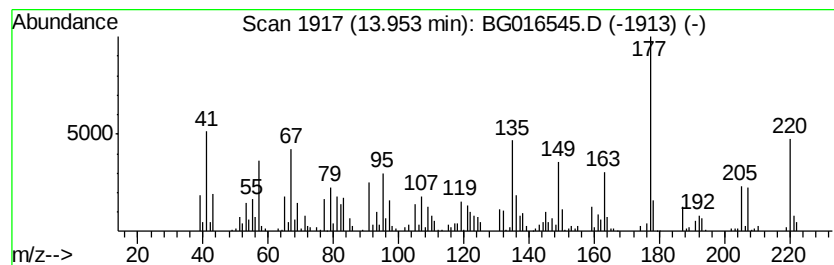
Instrument :
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG040615.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 2,5-Cyclohexadiene-1,4-dione... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.95	2.23 ng	132823	Acenaphthene-d10	14.30
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	2,5-Cyclohexadiene-1,4-dione, 2,...	220 C14H20O2	000719-22-2	96
2	Thiocoumarine, 4,4,6,7-tetramethyl-	220 C13H16OS	1000128-90-2	46
3	3-Buten-2-one, 4-(2,6,6-trimethy...	192 C13H20O	000079-77-6	38
4	N-[4-(Ethyl-methyl-amino)-phenyl...	192 C11H16N2O	099981-45-0	38
5	3-Buten-2-one, 4-(2,6,6-trimethy...	192 C13H20O	014901-07-6	35



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG041715\
Data File : BG016545.D
Acq On : 19 Apr 2015 5:30
Operator : TP/IZ
Sample : G1765-02
Misc :
ALS Vial : 54 Sample Multiplier: 1

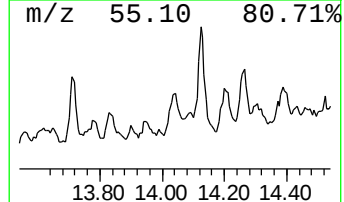
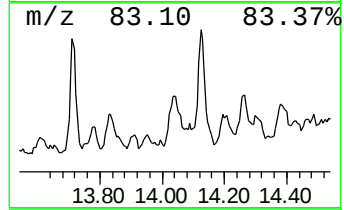
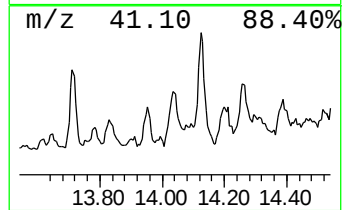
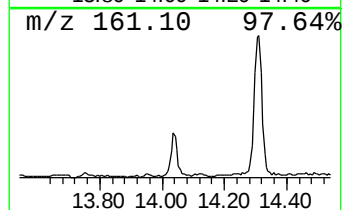
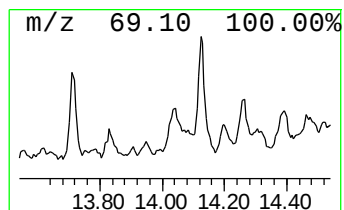
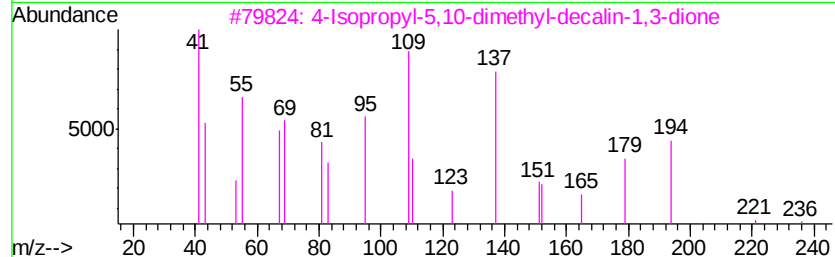
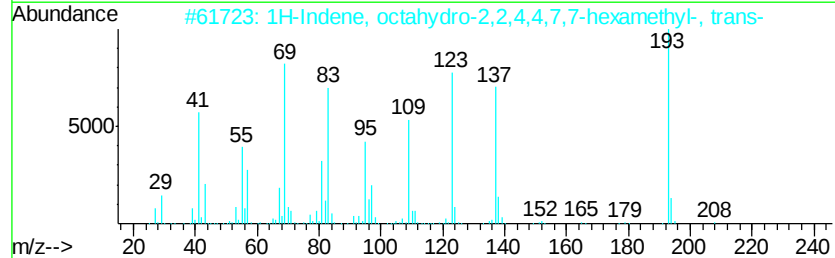
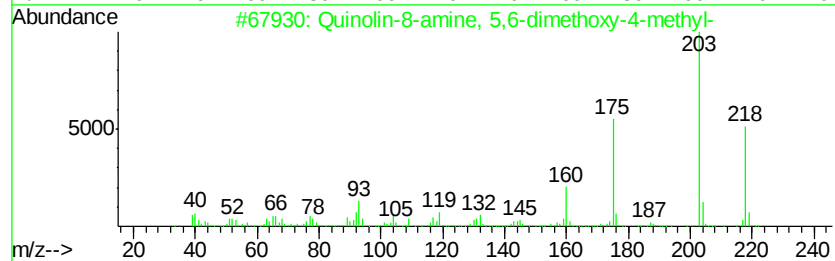
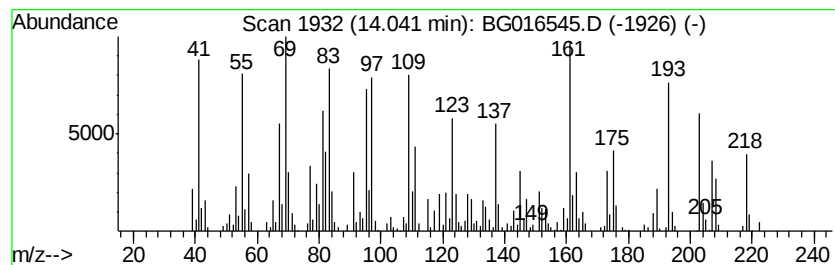
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG040615.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 unknown14.04 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.04	5.59 ng	332458	Acenaphthene-d10	14.30	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Quinolin-8-amine, 5,6-dimethoxy-...	218	C12H14N2O2	064992-95-6	15
2	1H-Indene, octahydro-2,2,4,4,7,7...	208	C15H28	054832-83-6	14
3	4-Isopropyl-5,10-dimethyl-decali...	236	C15H24O2	291278-94-9	14
4	6-Acetamido-1,4-benzodioxane	193	C10H11NO3	063546-19-0	14
5	2(1H)-Naphthalenone, octahydro-4...	194	C13H22O	054699-31-9	14



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG041715\
Data File : BG016545.D
Acq On : 19 Apr 2015 5:30
Operator : TP/IZ
Sample : G1765-02
Misc :
ALS Vial : 54 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
AS-2

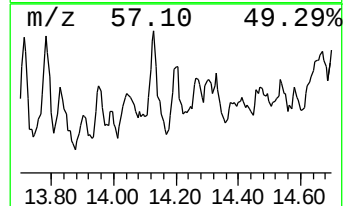
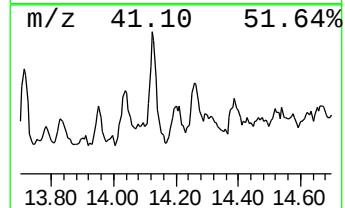
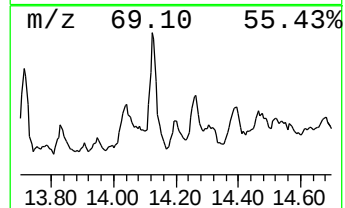
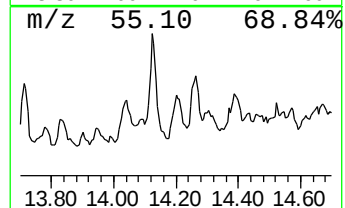
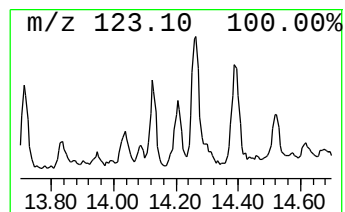
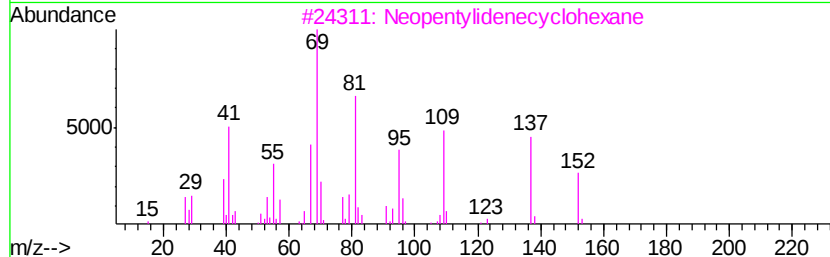
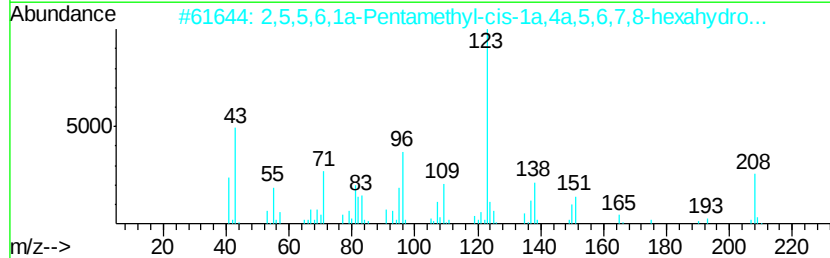
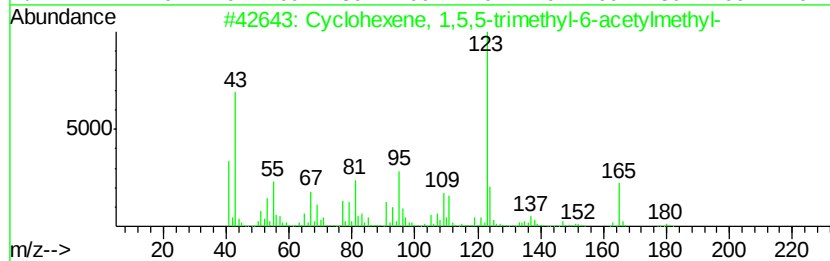
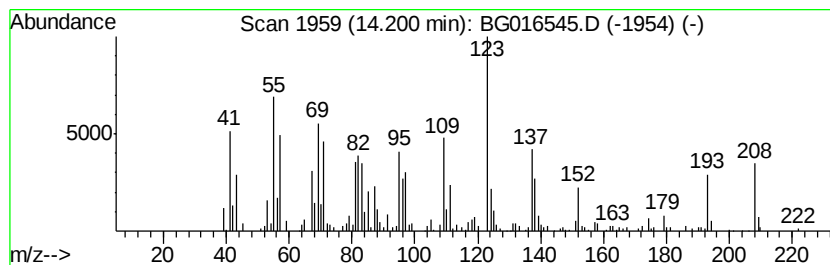
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 unknown14.20 Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.20	3.34 ng	198608	Acenaphthene-d10	14.30

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexene, 1,5,5-trimethyl-6-a...	180	C12H20O	211563-96-1	43
2		2,5,5,6,1a-Pentamethyl-cis-1a,4a...	208	C14H24O	1000215-77-9	38
3		Neopentylidenecyclohexane	152	C11H20	039546-80-0	38
4		1-Trimethylsilylpent-1-en-4-yne	138	C8H14Si	079516-25-9	35
5		o-Fluoroacetophenone	138	C8H7FO	000445-27-2	35



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG041715\
Data File : BG016545.D
Acq On : 19 Apr 2015 5:30
Operator : TP/IZ
Sample : G1765-02
Misc :
ALS Vial : 54 Sample Multiplier: 1

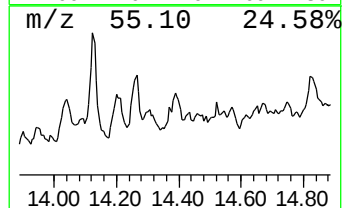
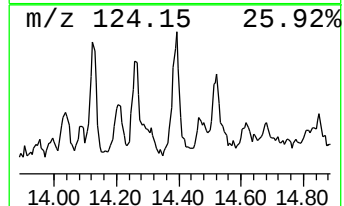
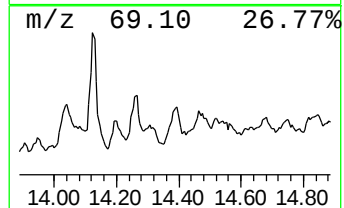
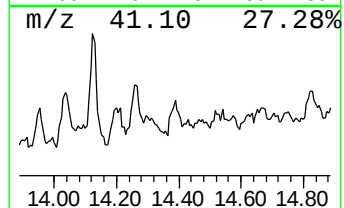
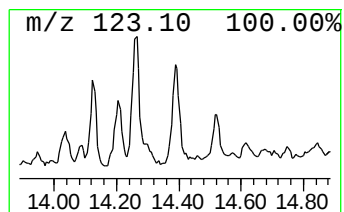
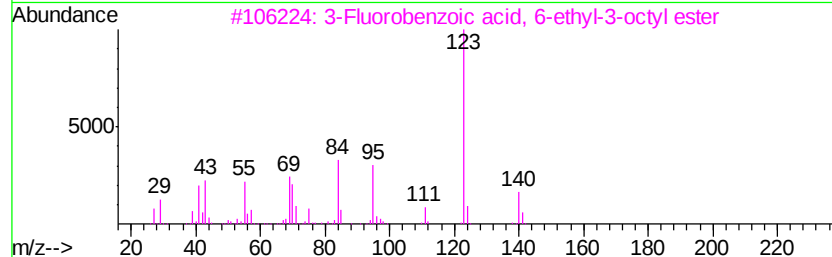
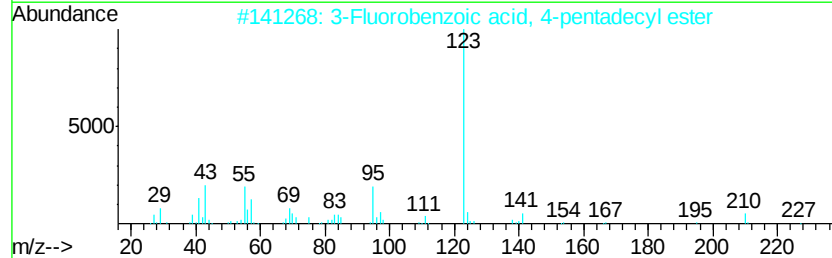
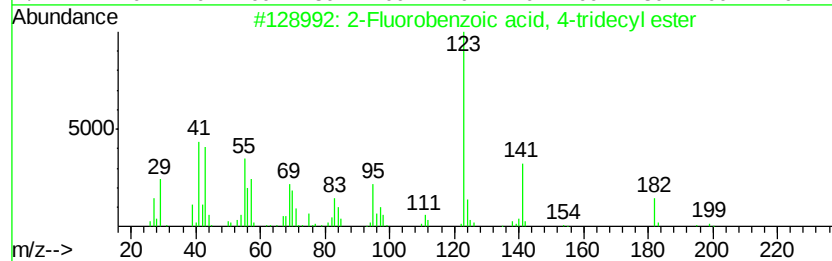
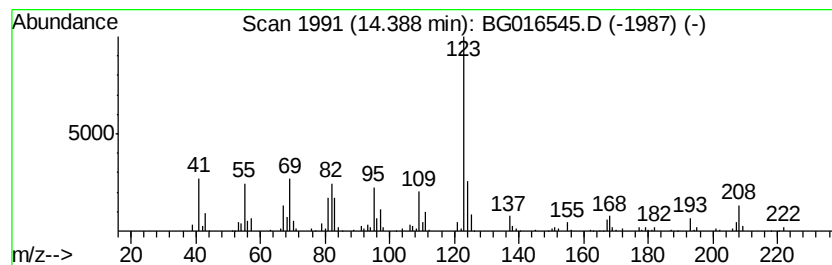
Instrument :
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ClientSampleId :
AS-2

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG040615.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 2-Fluorobenzoic acid, 4-tri... Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD		R.T.	
14.39	2.18 ng	129897	Acenaphthene-d10		14.30	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Fluorobenzoic acid,	4-tridecyl...	322	C20H31F02	1000280-61-2	50
2	3-Fluorobenzoic acid,	4-pentadec...	350	C22H35F02	1000280-60-9	47
3	3-Fluorobenzoic acid,	6-ethyl-3-...	280	C17H25F02	1000279-01-9	47
4	2-Fluorobenzoic acid,	6-ethyl-2-...	280	C17H25F02	1000278-96-7	47
5	4-Fluorobenzoic acid,	3-pentadec...	350	C22H35F02	1000280-68-4	47



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG041715\
Data File : BG016545.D
Acq On : 19 Apr 2015 5:30
Operator : TP/IZ
Sample : G1765-02
Misc :
ALS Vial : 54 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
AS-2

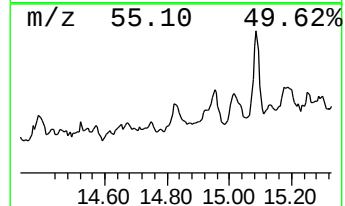
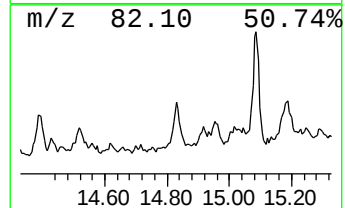
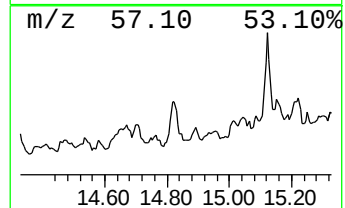
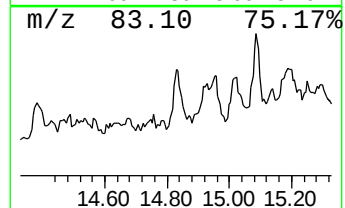
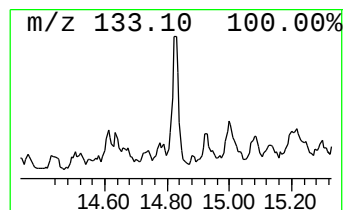
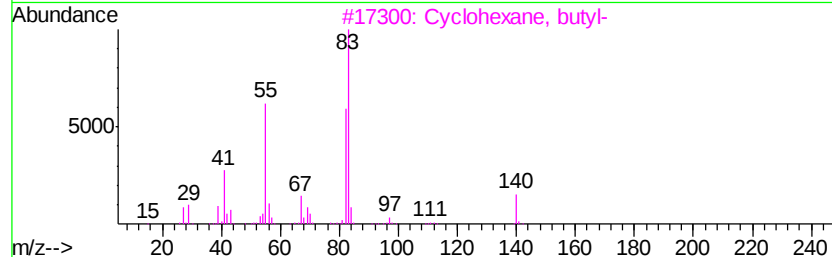
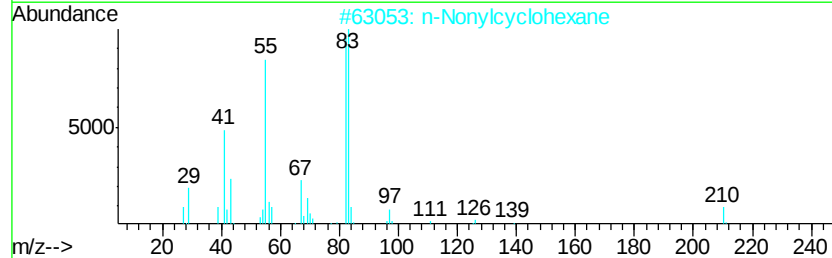
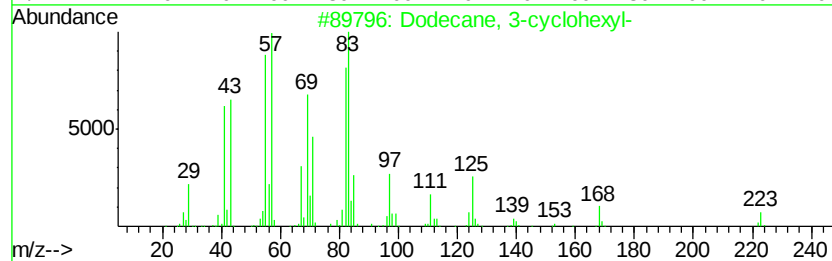
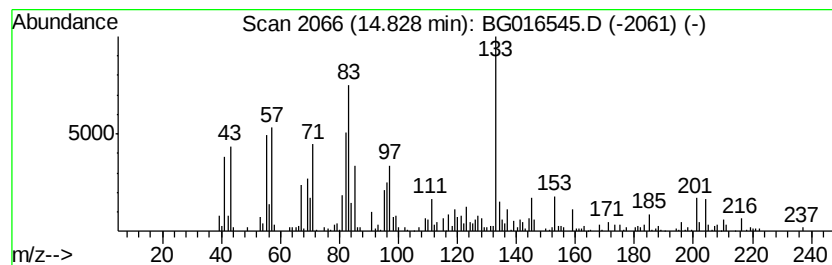
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 unknown14.83 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.83	4.15 ng	247193	Acenaphthene-d10	14.30

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Dodecane, 3-cyclohexyl-	252	C18H36	013151-83-2	35
2		n-Nonylcyclohexane	210	C15H30	002883-02-5	30
3		Cyclohexane, butyl-	140	C10H20	001678-93-9	22
4		Cyclohexane, octyl-	196	C14H28	001795-15-9	22
5		Pentadecanal-	226	C15H30O	002765-11-9	22



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG041715\
Data File : BG016545.D
Acq On : 19 Apr 2015 5:30
Operator : TP/IZ
Sample : G1765-02
Misc :
ALS Vial : 54 Sample Multiplier: 1

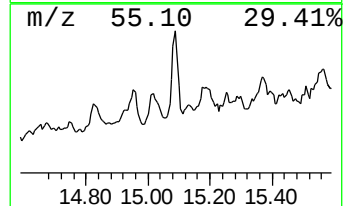
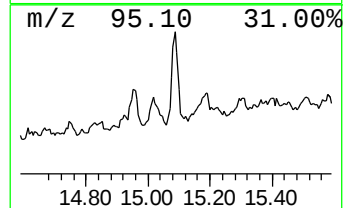
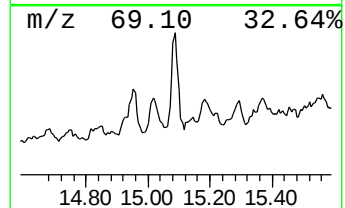
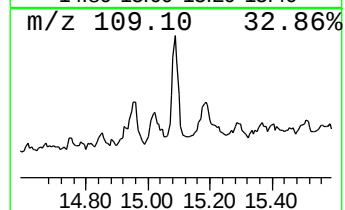
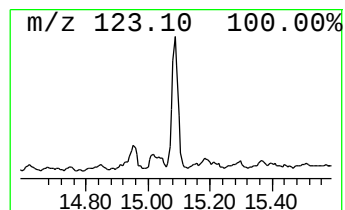
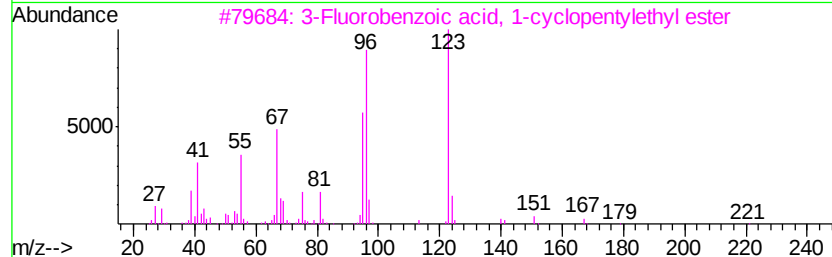
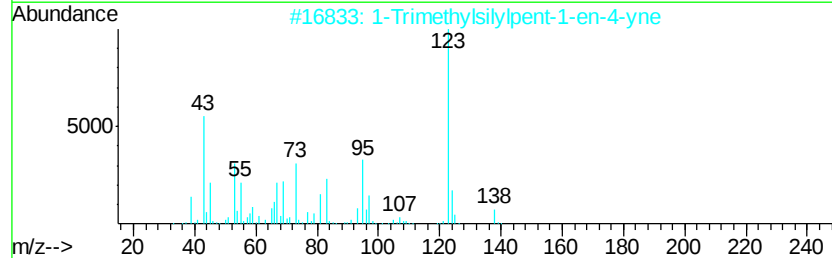
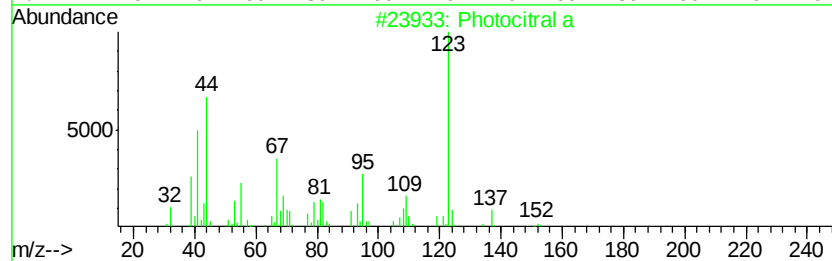
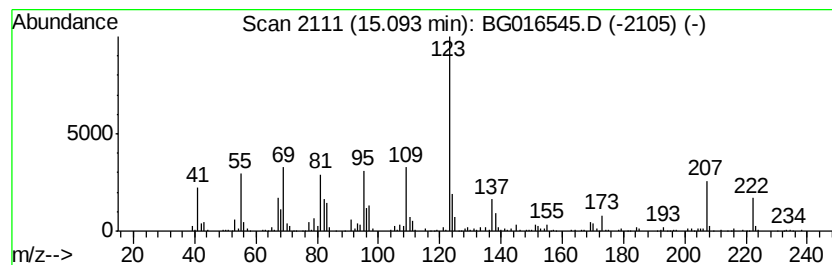
Instrument :
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ClientSampleId :
AS-2

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG040615.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 unknown15.09 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD		R.T.	
15.09	9.79 ng	582552	Acenaphthene-d10		14.30	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Photocitral a	152	C10H16O	1000292-85-0	47
2		1-Trimethylsilylpent-1-en-4-yne	138	C8H14Si	079516-25-9	47
3		3-Fluorobenzoic acid, 1-cyclopent...	236	C14H17FO2	1000279-04-7	47
4		1-Butanone, 1-bicyclo[4.1.0]hept...	166	C11H18O	054764-62-4	43
5		2-(2-Hydroxycyclohexyl)-furan	166	C10H14O2	1000145-92-8	38



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG041715\
Data File : BG016545.D
Acq On : 19 Apr 2015 5:30
Operator : TP/IZ
Sample : G1765-02
Misc :
ALS Vial : 54 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
AS-2

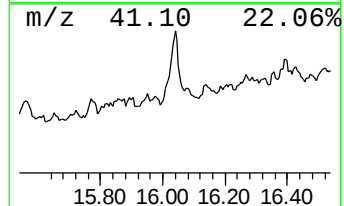
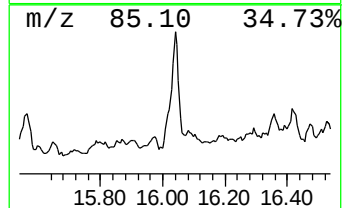
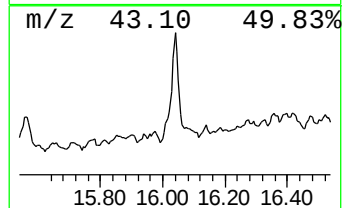
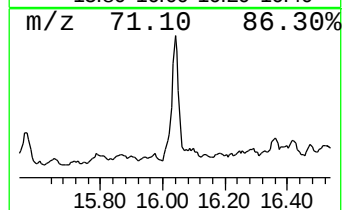
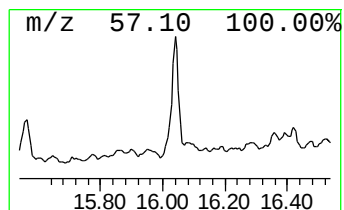
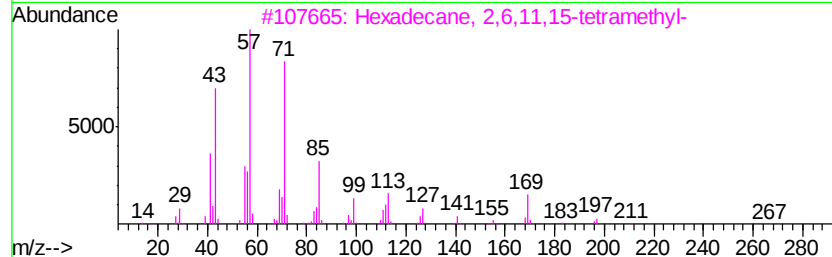
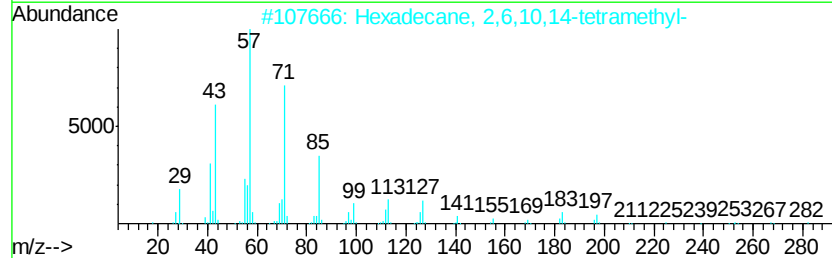
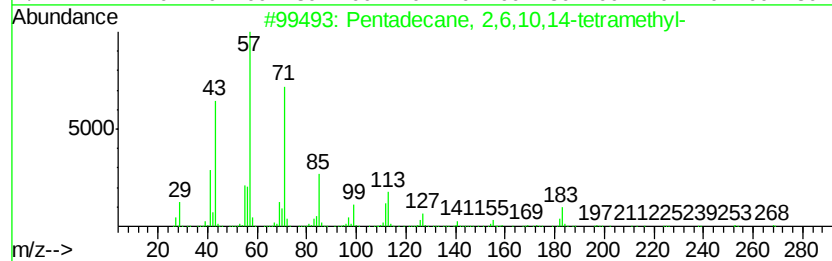
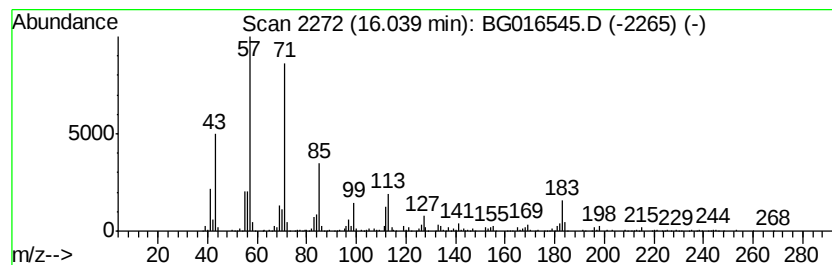
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 Pentadecane, 2,6,10,14-tetr... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.04	18.01 ng	1015590	Phenanthrene-d10	17.04

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentadecane, 2,6,10,14-tetramethyl-	268	C19H40	001921-70-6	81
2		Hexadecane, 2,6,10,14-tetramethyl-	282	C20H42	000638-36-8	80
3		Hexadecane, 2,6,11,15-tetramethyl-	282	C20H42	000504-44-9	80
4		Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	64
5		Hexadecane, 2,6,10-trimethyl-	268	C19H40	055000-52-7	64



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG041715\
Data File : BG016545.D
Acq On : 19 Apr 2015 5:30
Operator : TP/IZ
Sample : G1765-02
Misc :
ALS Vial : 54 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
AS-2

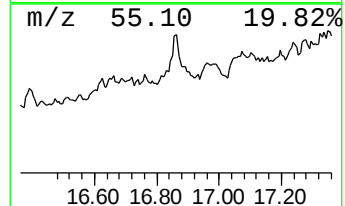
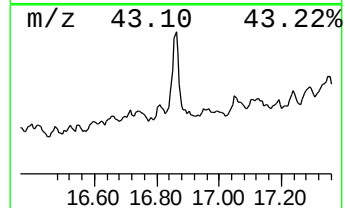
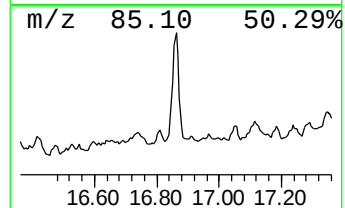
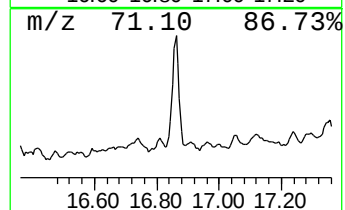
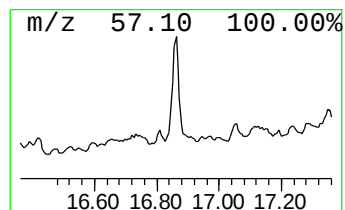
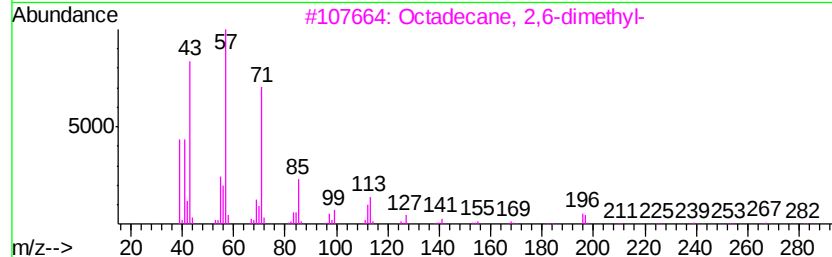
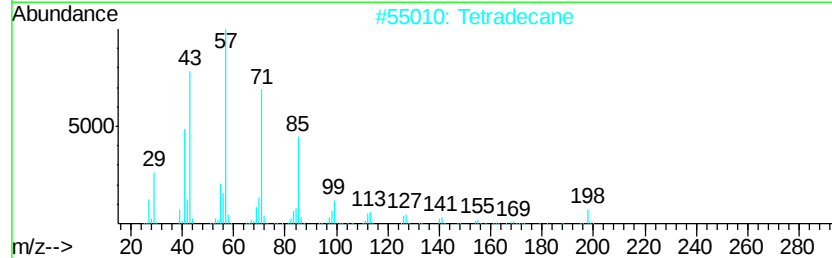
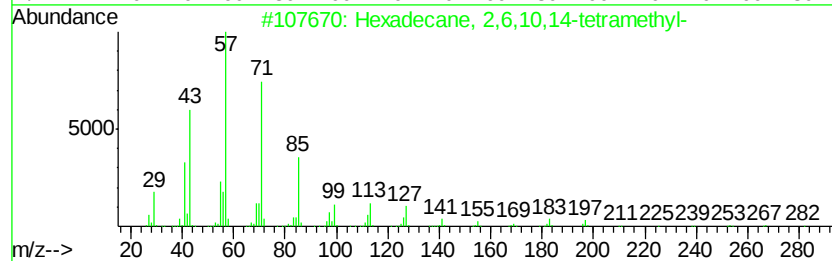
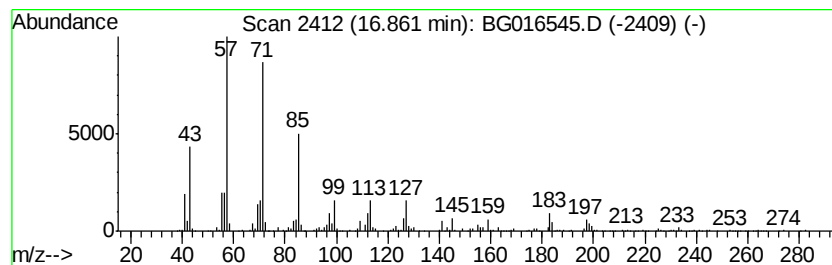
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG040615.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 Hexadecane, 2,6,10,14-tetra... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.86	14.96 ng	843206	Phenanthrene-d10	17.04

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecane, 2,6,10,14-tetramethyl-	282	C20H42	000638-36-8	99
2		Tetradecane	198	C14H30	000629-59-4	90
3		Octadecane, 2,6-dimethyl-	282	C20H42	075163-97-2	86
4		Heptacosane	380	C27H56	000593-49-7	80
5		Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	80



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG041715\
 Data File : BG016545.D
 Acq On : 19 Apr 2015 5:30
 Operator : TP/IZ
 Sample : G1765-02
 Misc :
 ALS Vial : 54 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 AS-2

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG040615.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...	2.83	6.5	ng	169835	1	7.65	522019	20.0
Hexane, 2,2-dimet...	2.92	12.3	ng	321335	1	7.65	522019	20.0
2-Pentanone, 4-hy...	4.77	15.8	ng	411065	1	7.65	522019	20.0
unknown7.20	7.20	50.3	ng	1312880	1	7.65	522019	20.0
Decahydro-4,4,8,9...	13.71	4.6	ng	272931	3	14.30	1189910	20.0
1H-Indene, octahy...	13.83	2.4	ng	139959	3	14.30	1189910	20.0
2,5-Cyclohexadien...	13.95	2.2	ng	132823	3	14.30	1189910	20.0
unknown14.04	14.04	5.6	ng	332458	3	14.30	1189910	20.0
unknown14.20	14.20	3.3	ng	198608	3	14.30	1189910	20.0
2-Fluorobenzoic a...	14.39	2.2	ng	129897	3	14.30	1189910	20.0
unknown14.83	14.83	4.2	ng	247193	3	14.30	1189910	20.0
unknown15.09	15.09	9.8	ng	582552	3	14.30	1189910	20.0
Pentadecane, 2,6,...	16.04	18.0	ng	1015590	4	17.04	1127500	20.0
Hexadecane, 2,6,1...	16.86	15.0	ng	843206	4	17.04	1127500	20.0