

Data Path : Z:\HPCHEM1\BNA G\DATA\BG012518\
 Data File : BG032317.D
 Acq On : 26 Jan 2018 6:39
 Operator : SJ/JU
 Sample : J1267-03
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 IGW-6

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-BG011218.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	3.466	24	30	44	rVB	281659	422600	14.13%	2.572%
2	4.630	221	228	238	rVB	45373	75428	2.52%	0.459%
3	5.652	391	402	415	rBV	533996	885692	29.62%	5.390%
4	6.087	470	476	482	rBV	99058	164668	5.51%	1.002%
5	6.157	482	488	494	rVV	291429	509563	17.04%	3.101%
6	6.228	494	500	514	rVB	203722	483896	16.18%	2.945%
7	6.627	561	568	575	rBV	128099	219563	7.34%	1.336%
8	7.650	736	742	748	rBV2	21973	36744	1.23%	0.224%
9	7.761	755	761	766	rBV	80710	138590	4.63%	0.843%
10	7.832	766	773	780	rVV2	318440	593456	19.85%	3.611%
11	7.902	780	785	794	rVB3	56134	94880	3.17%	0.577%
12	8.079	808	815	822	rBB	47458	81663	2.73%	0.497%
13	8.237	834	842	850	rBV	361162	677182	22.65%	4.121%
14	8.349	854	861	871	rVB	210815	363536	12.16%	2.212%
15	8.725	918	925	936	rVB2	95403	182474	6.10%	1.110%
16	8.837	936	944	951	rBV	81498	143163	4.79%	0.871%
17	9.060	974	982	987	rBV	390677	724969	24.24%	4.412%
18	9.107	987	990	998	rVB	58669	98884	3.31%	0.602%
19	9.277	1013	1019	1027	rVB4	17411	30965	1.04%	0.188%
20	9.365	1027	1034	1042	rBV2	28552	52646	1.76%	0.320%
21	9.853	1111	1117	1121	rBV3	27537	50771	1.70%	0.309%
22	9.918	1121	1128	1141	rVB	289734	573950	19.19%	3.493%
23	10.447	1212	1218	1225	rVB	26683	44101	1.47%	0.268%
24	10.823	1276	1282	1290	rBV2	22818	40788	1.36%	0.248%
25	10.975	1301	1308	1316	rBV3	56981	112632	3.77%	0.685%
26	11.604	1405	1415	1419	rBV	130769	250390	8.37%	1.524%
27	11.651	1419	1423	1430	rVB	119449	217958	7.29%	1.326%
28	13.220	1682	1690	1698	rVB	85709	138611	4.64%	0.843%
29	13.431	1716	1726	1735	rBV	49553	95629	3.20%	0.582%
30	13.984	1812	1820	1833	rBV	1131851	1825828	61.06%	11.111%
31	15.358	2048	2054	2063	rVB	253359	425379	14.22%	2.589%
32	16.686	2275	2280	2288	rVB	59525	86379	2.89%	0.526%
33	16.833	2298	2305	2313	rBV	783232	1227830	41.06%	7.472%
34	18.090	2512	2519	2532	rBV	412235	644123	21.54%	3.920%

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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 >
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35	20.629	2944	2951	2962	rVB	2206440	2990364	100.00%	18.197%
36	22.421	3245	3256	3263	rBV	457896	851751	28.48%	5.183%
37	26.099	3869	3882	3895	rBV2	275699	875951	29.29%	5.330%

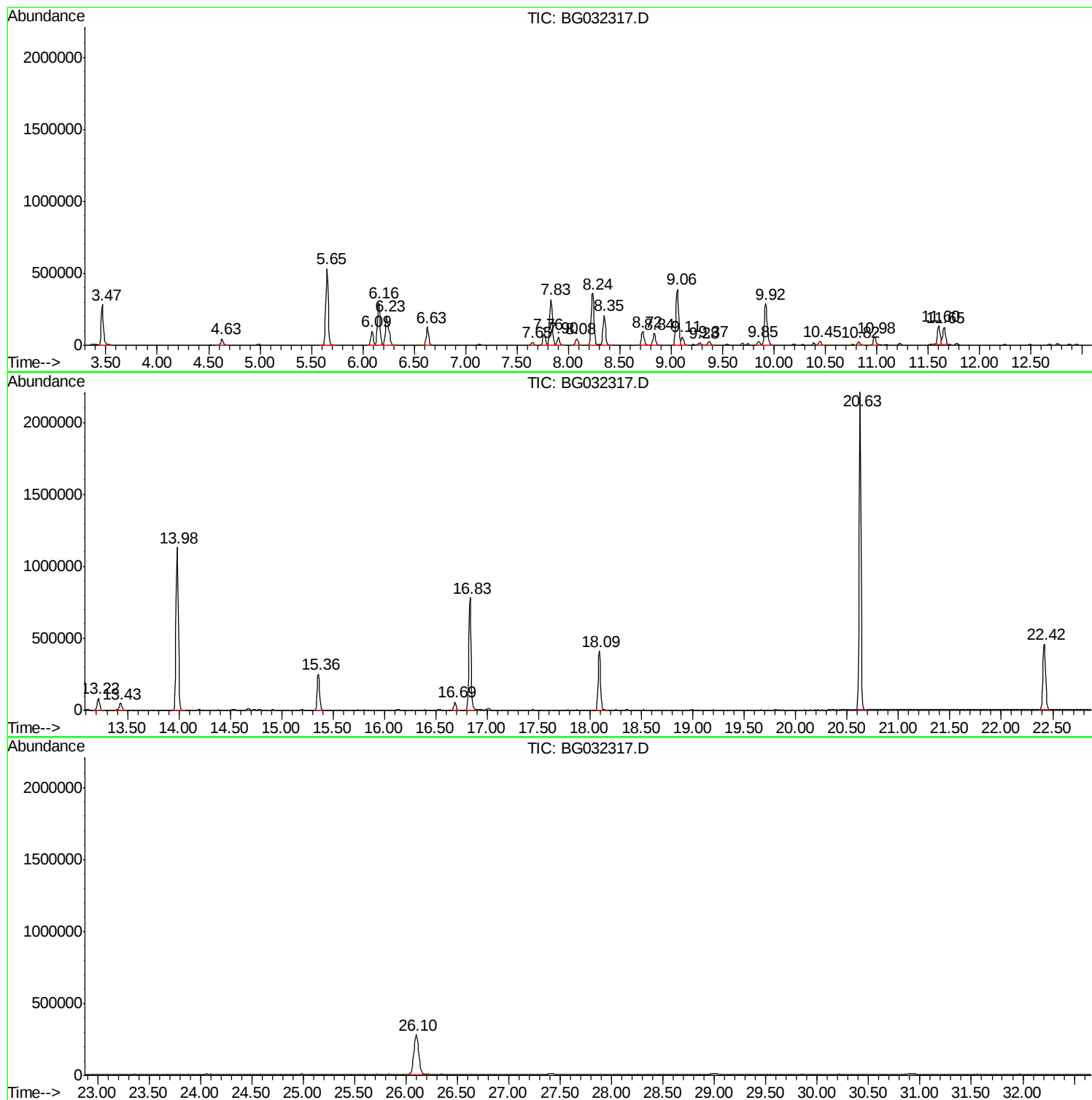
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Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG011218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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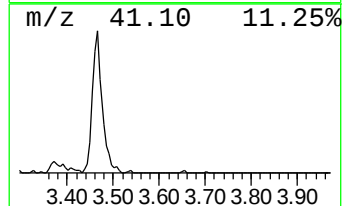
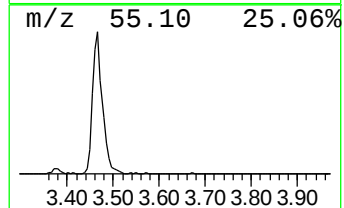
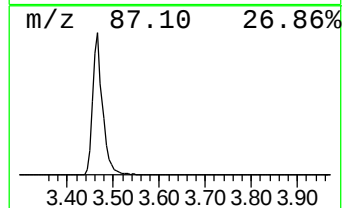
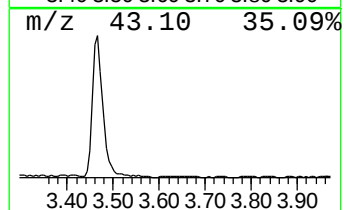
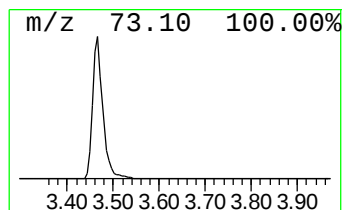
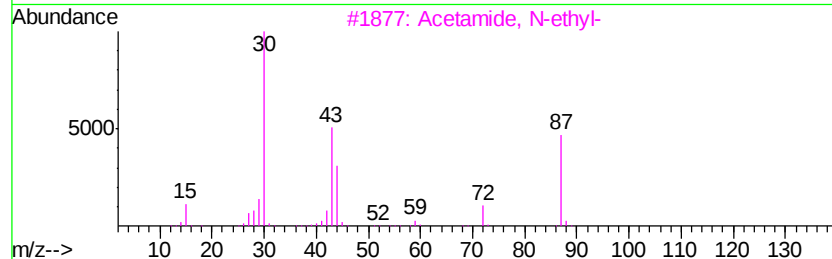
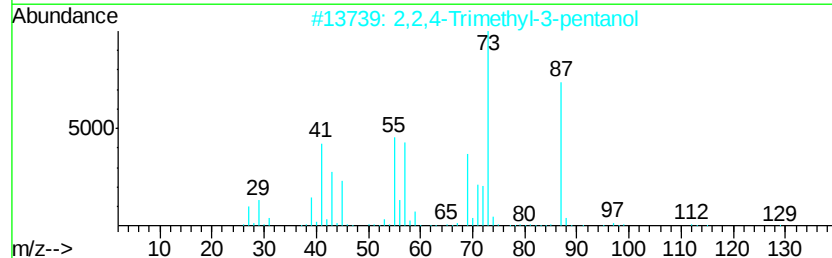
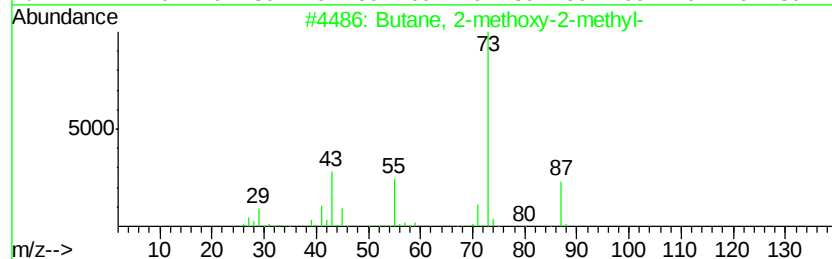
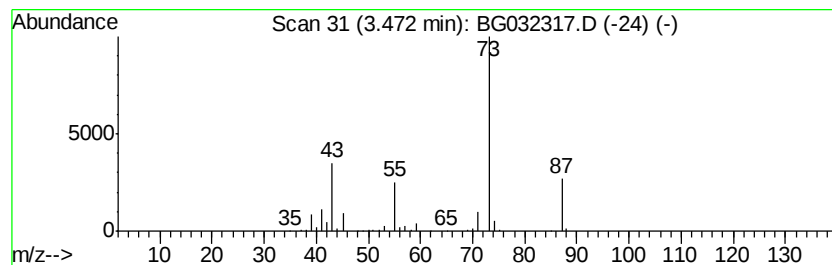
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.47	46.32 ng	422600	1,4-Dichlorobenzene-d4	8.72

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	38
3		Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	22
4		N-Ethylformamide	73	C3H7NO	000627-45-2	9
5		Silane, ethyltrimethyl-	102	C5H14Si	003439-38-1	9



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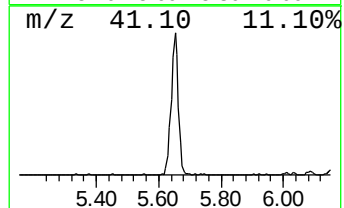
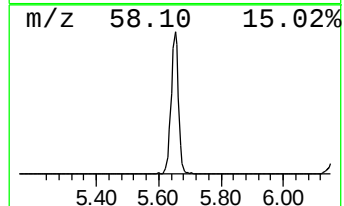
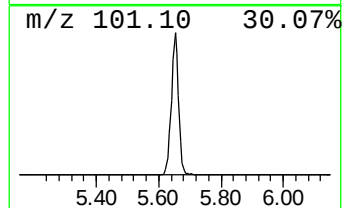
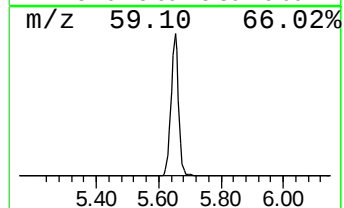
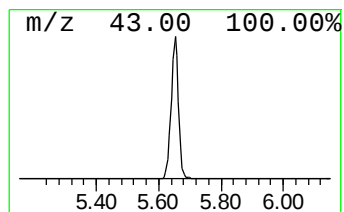
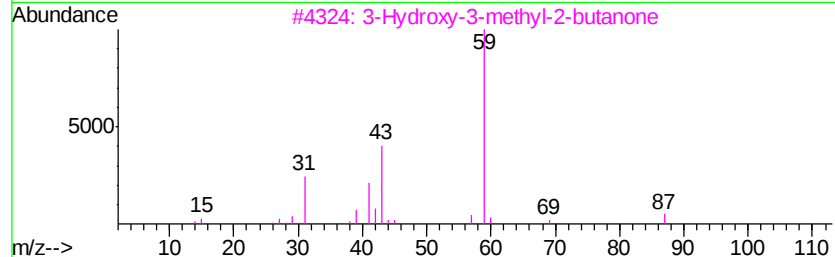
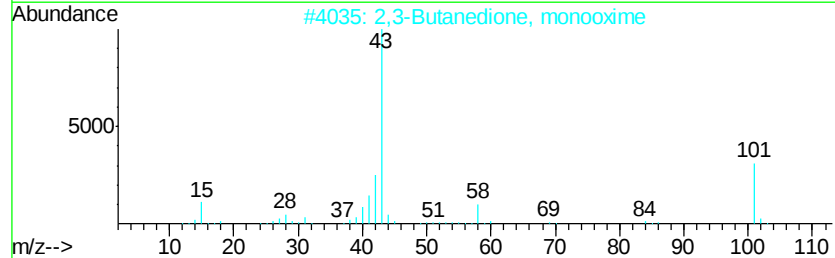
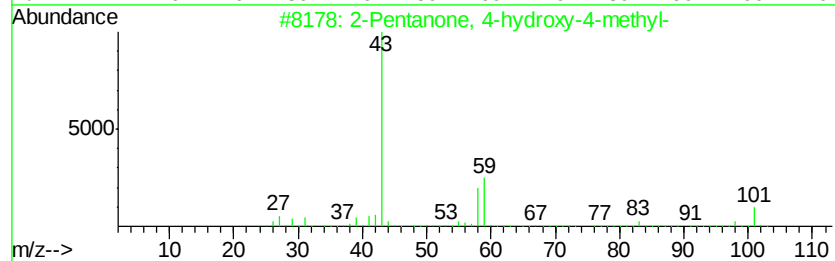
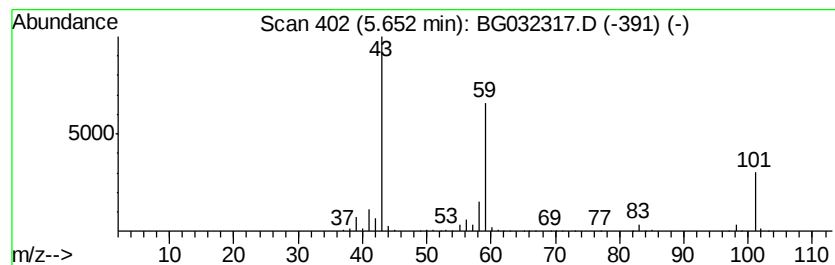
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TIC Library : C:\Database\NIST11.L
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Peak Number 3 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.65	97.08 ng	885692	1,4-Dichlorobenzene-d4	8.72

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	25
3		3-Hydroxy-3-methyl-2-butanone	102	C5H10O2	000115-22-0	16
4		5-Hexen-2-one	98	C6H10O	000109-49-9	9
5		Butane	58	C4H10	000106-97-8	9



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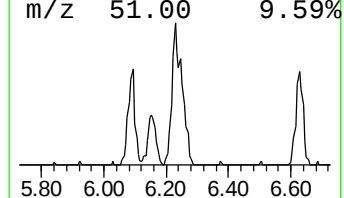
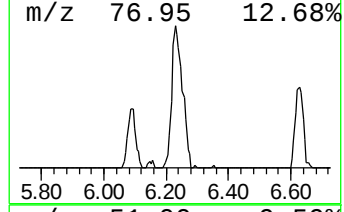
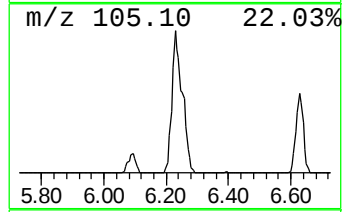
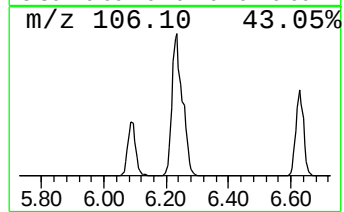
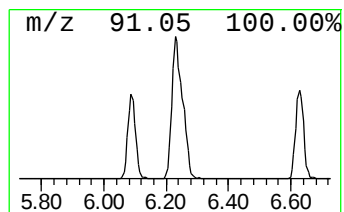
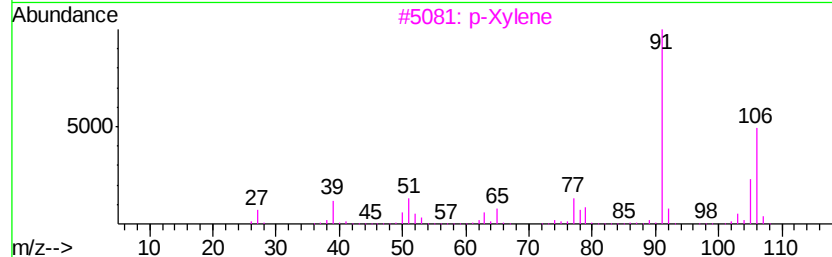
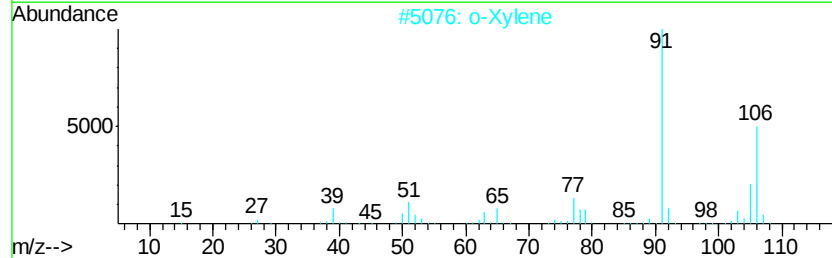
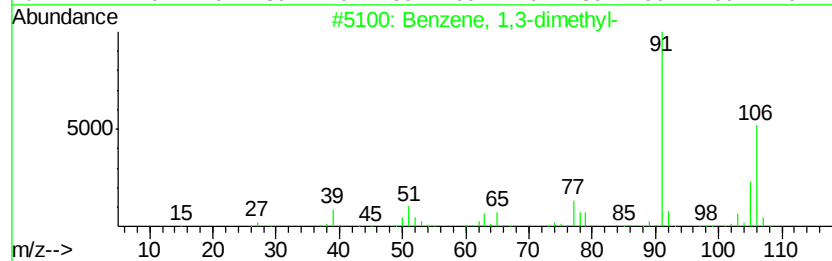
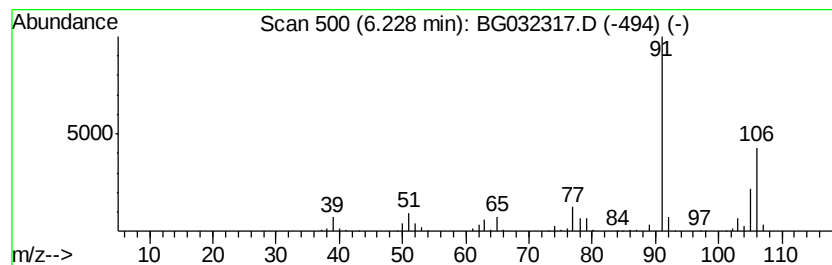
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Benzene, 1,3-dimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.23	53.04 ng	483896	1,4-Dichlorobenzene-d4	8.72

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	97
2		o-Xylene	106	C8H10	000095-47-6	97
3		p-Xylene	106	C8H10	000106-42-3	97
4		Ethylbenzene	106	C8H10	000100-41-4	74
5		1,3-Cyclopentadiene, 5-(1-methyl...	106	C8H10	002175-91-9	74



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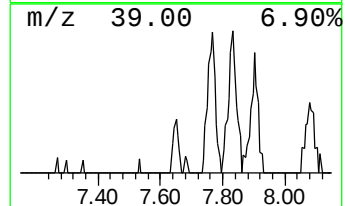
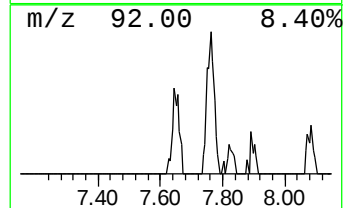
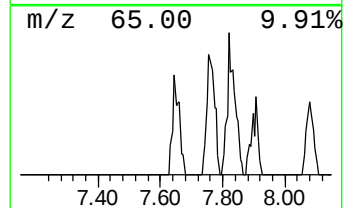
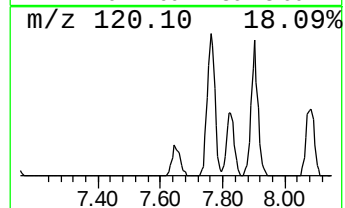
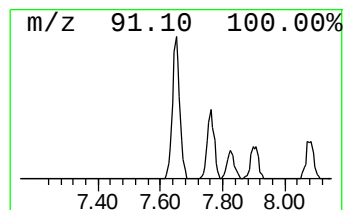
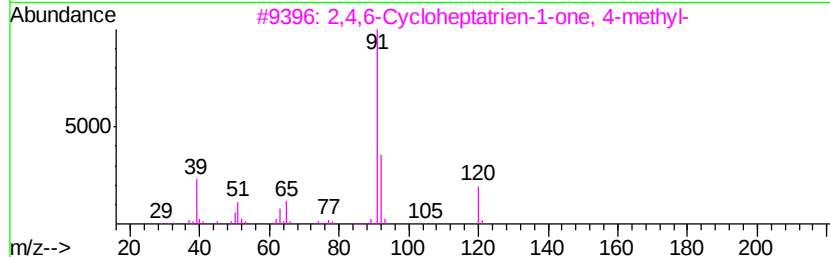
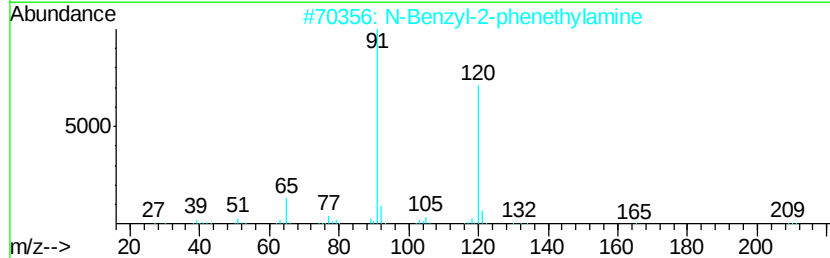
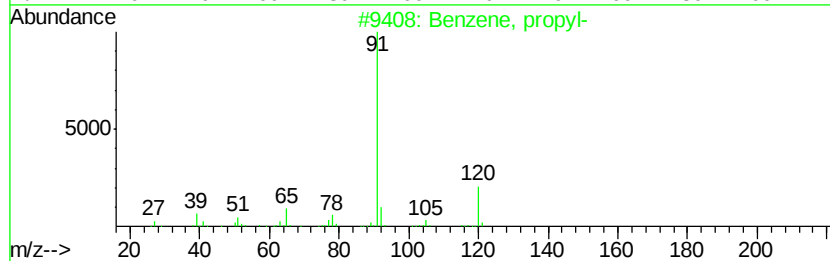
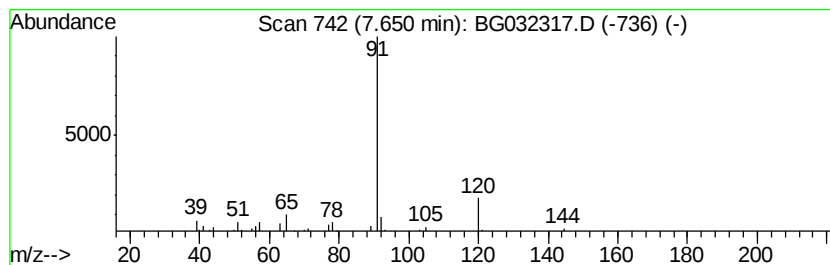
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Benzene, propyl- Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.65	4.03 ng	36744	1,4-Dichlorobenzene-d4	8.72

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, propyl-	120	C9H12	000103-65-1	87
2		N-Benzyl-2-phenethylamine	211	C15H17N	003647-71-0	78
3		2,4,6-Cycloheptatrien-1-one, 4-m...	120	C8H8O	1000327-34-9	53
4		Acetyl benzyl disulfide	198	C9H10S2	005797-02-4	47
5		Hydrazinecarboxylic acid, phenyl...	166	C8H10N2O2	005331-43-1	47



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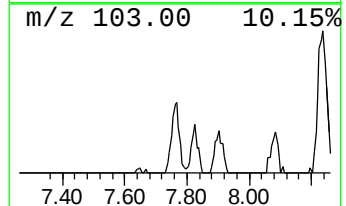
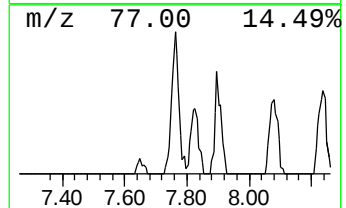
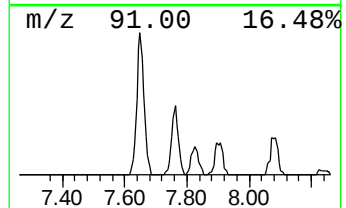
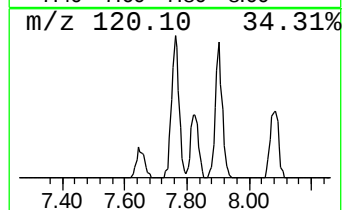
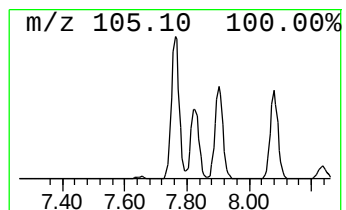
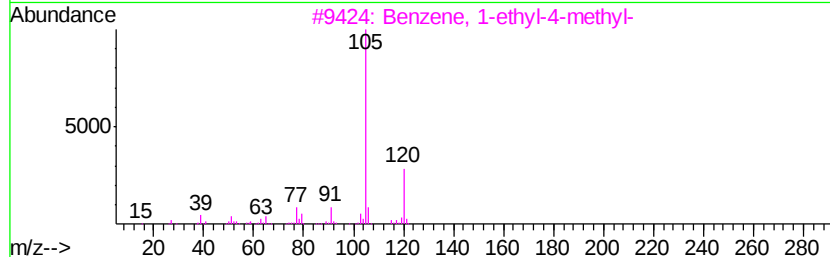
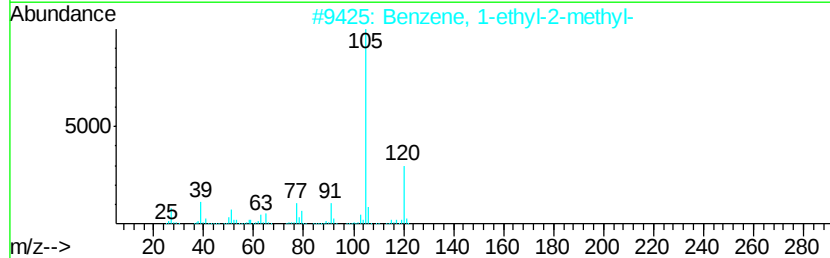
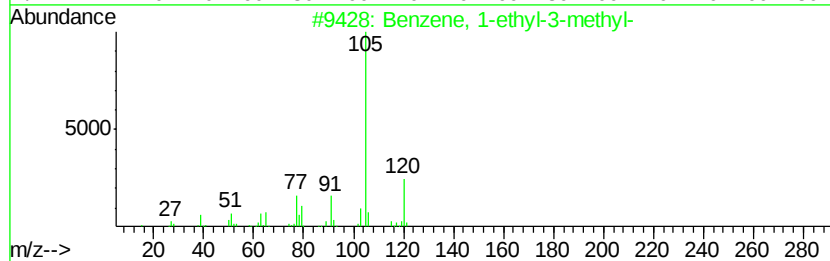
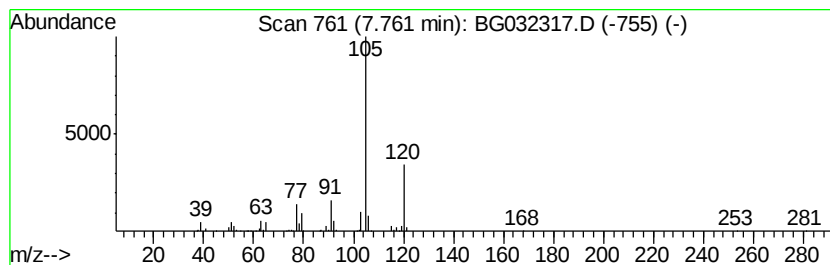
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Peak Number 8 Benzene, 1-ethyl-3-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.76	15.19 ng	138590	1,4-Dichlorobenzene-d4	8.72

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
2		Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	91
3		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
4		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	87
5		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	87



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Sample : J1267-03
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
IGW-6

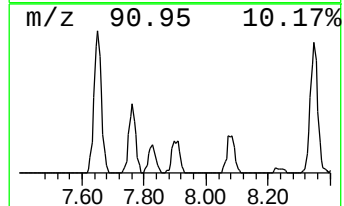
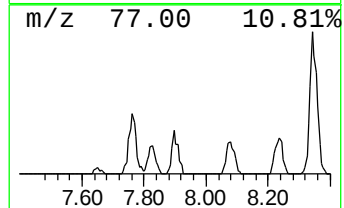
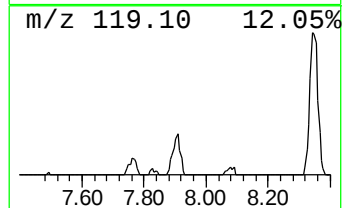
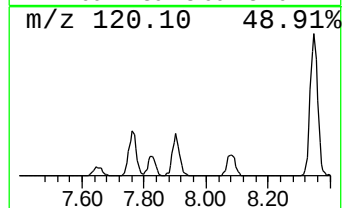
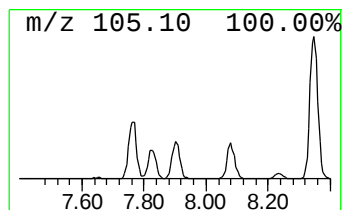
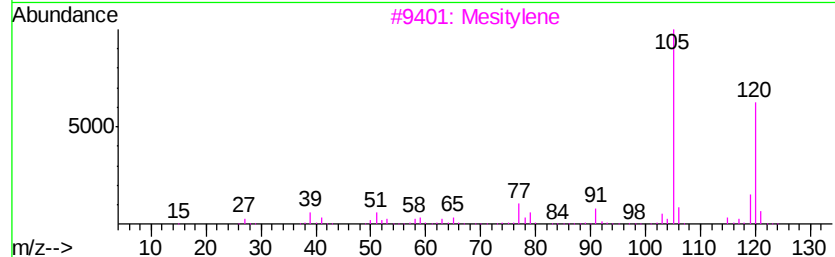
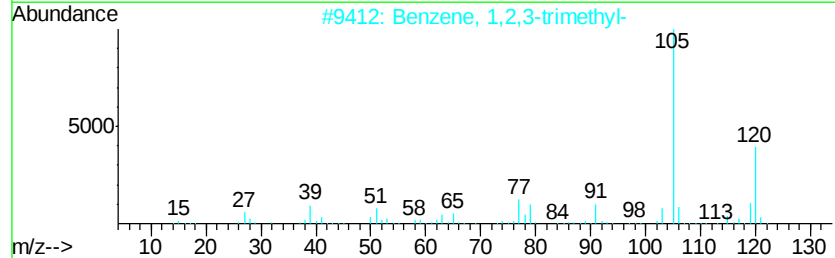
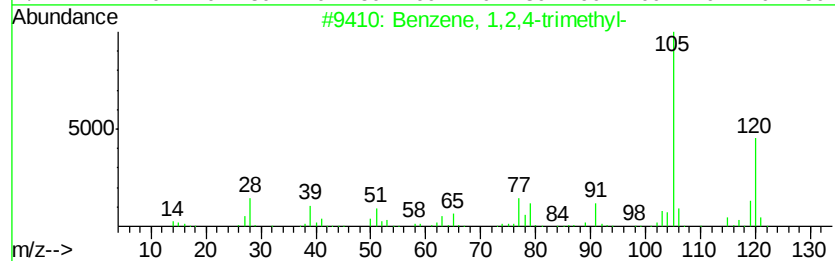
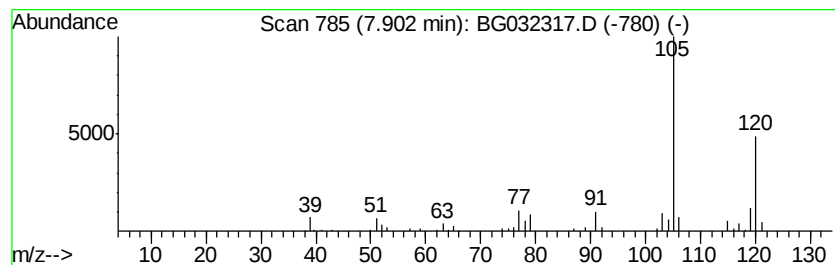
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Benzene, 1,2,4-trimethyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.90	10.40 ng	94880	1,4-Dichlorobenzene-d4	8.72

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	96
2		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
3		Mesitylene	120	C9H12	000108-67-8	91
4		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	90
5		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	87



Data Path : Z:\HPCHEM1\BNA G\DATA\BG012518\
Data File : BG032317.D
Acq On : 26 Jan 2018 6:39
Operator : SJ/JU
Sample : J1267-03
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
IGW-6

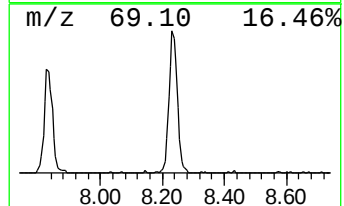
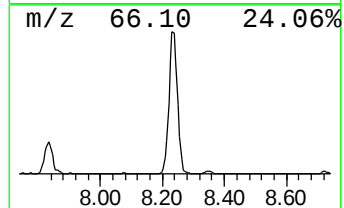
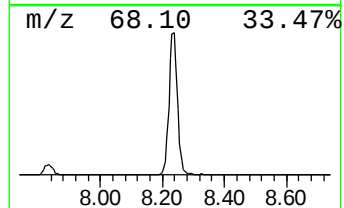
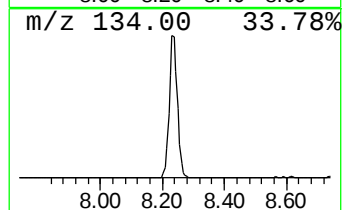
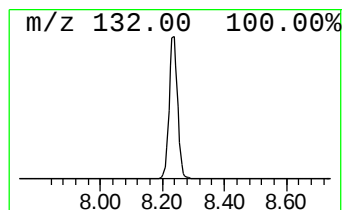
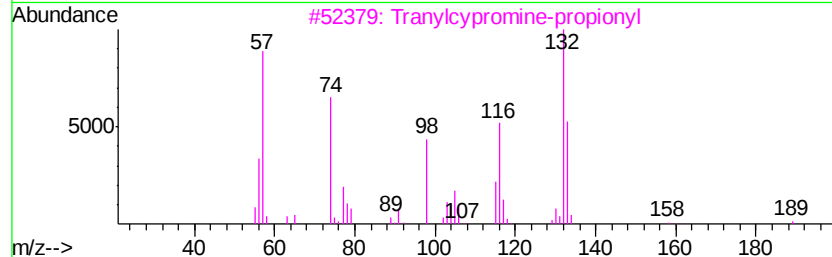
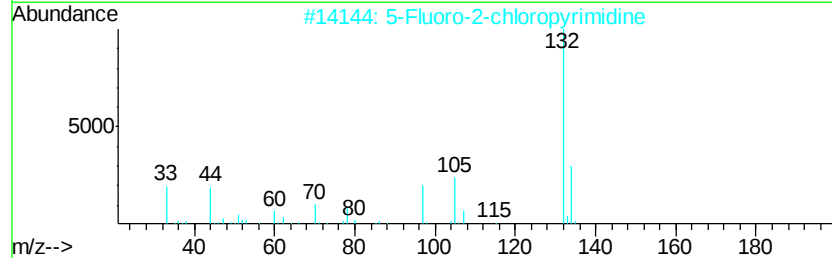
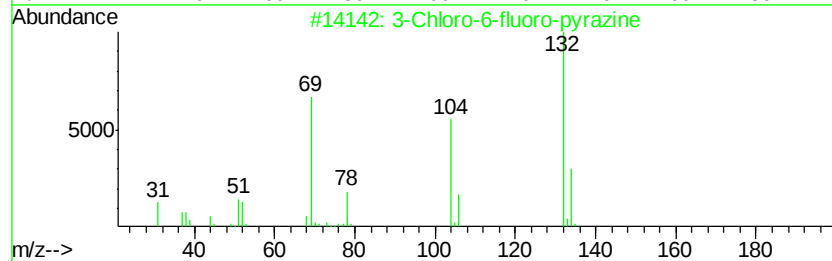
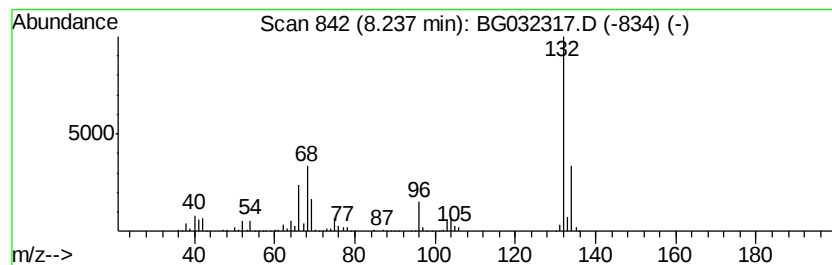
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 unknown8.24 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.24	74.22 ng	677182	1,4-Dichlorobenzene-d4	8.72

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	35
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	17
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	12
4		1,3,5-Trifluorobenzene	132	C6H3F3	000372-38-3	10
5		N-(-3,4-Methylenedioxyphenylisop...	267	C17H17NO2	1000378-96-6	10



Data Path : Z:\HPCHEM1\BNA G\DATA\BG012518\
Data File : BG032317.D
Acq On : 26 Jan 2018 6:39
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Instrument :
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ClientSampleId :
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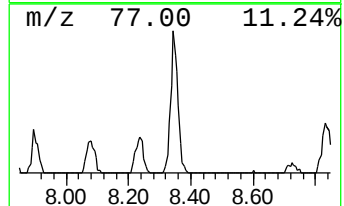
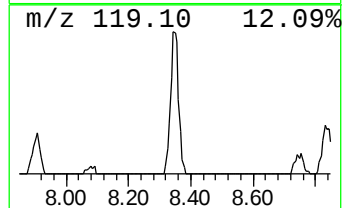
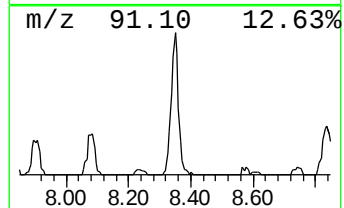
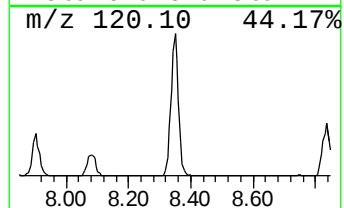
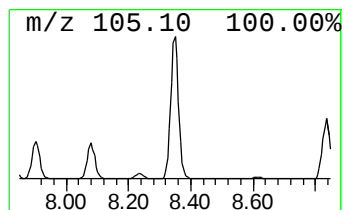
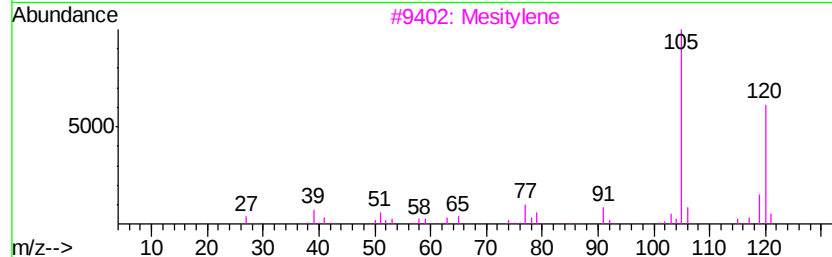
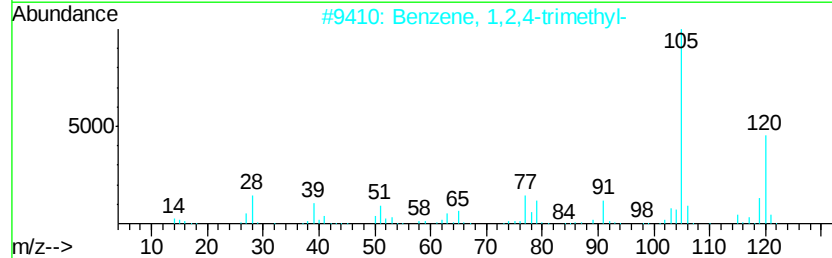
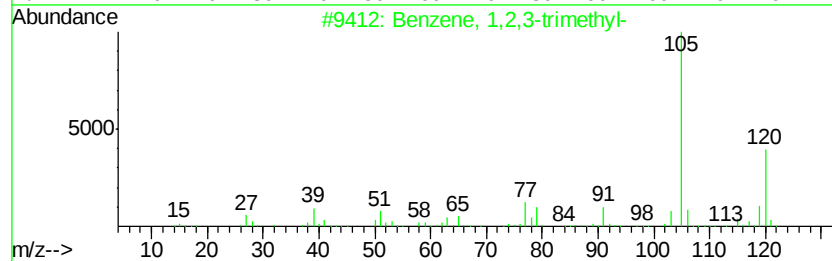
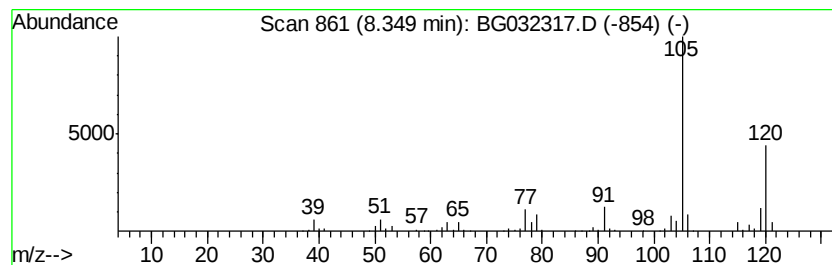
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Benzene, 1,2,3-trimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.35	39.85 ng	363536	1,4-Dichlorobenzene-d4	8.72

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	97
2		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
3		Mesitylene	120	C9H12	000108-67-8	91
4		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
5		Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	90



Data Path : Z:\HPCHEM1\BNA G\DATA\BG012518\
Data File : BG032317.D
Acq On : 26 Jan 2018 6:39
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Sample : J1267-03
Misc :
ALS Vial : 22 Sample Multiplier: 1

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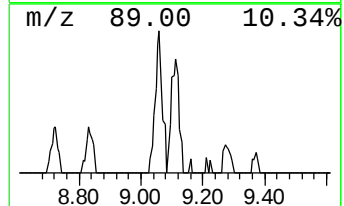
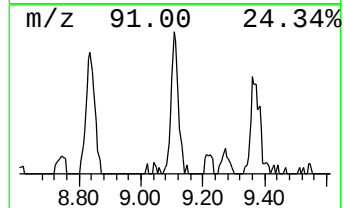
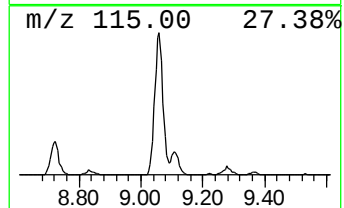
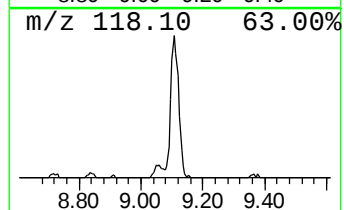
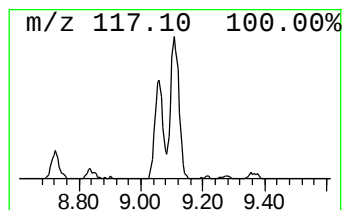
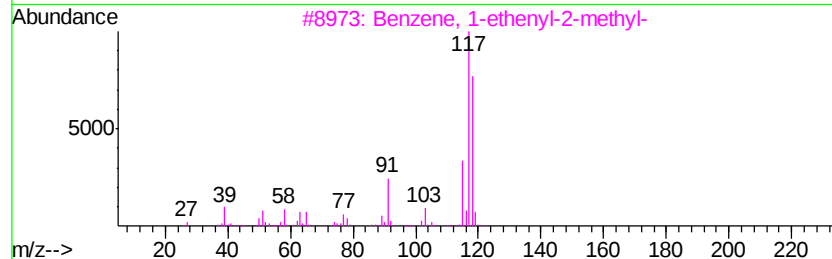
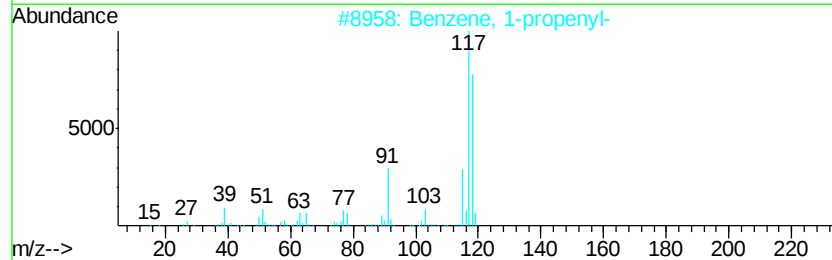
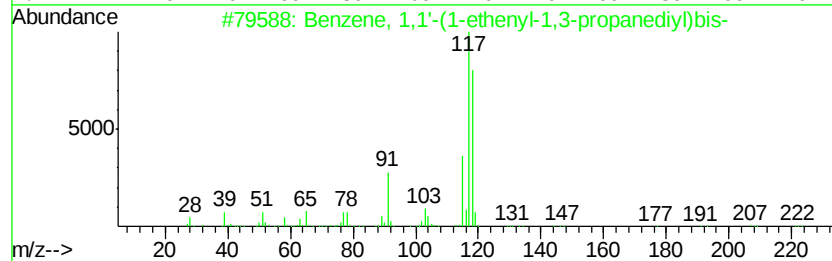
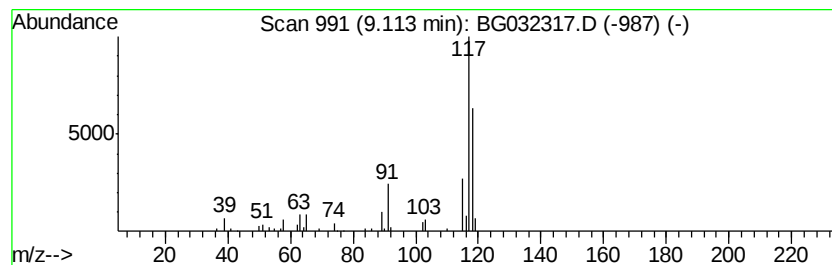
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Benzene, 1,1'-(1-ethenyl-1,... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.11	10.84 ng	98884	1,4-Dichlorobenzene-d4	8.72

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,1'-(1-ethenyl-1,3-pro...	222	C17H18	061141-97-7	90
2		Benzene, 1-propenyl-	118	C9H10	000637-50-3	87
3		Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	87
4		Indane	118	C9H10	000496-11-7	83
5		Benzene, 2-propenyl-	118	C9H10	000300-57-2	74



Data Path : Z:\HPCHEM1\BNA G\DATA\BG012518\
Data File : BG032317.D
Acq On : 26 Jan 2018 6:39
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Sample : J1267-03
Misc :
ALS Vial : 22 Sample Multiplier: 1

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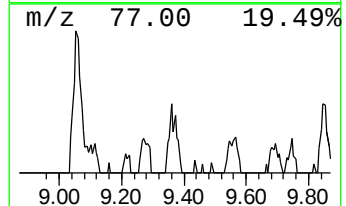
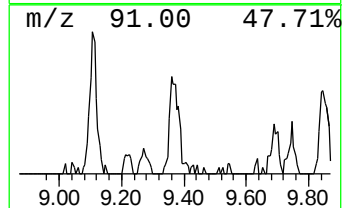
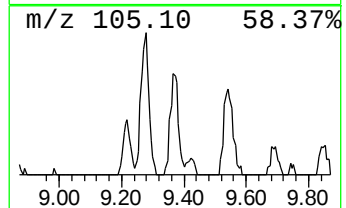
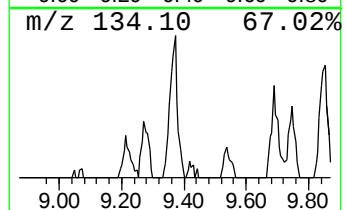
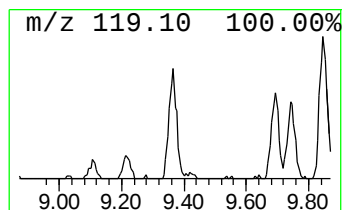
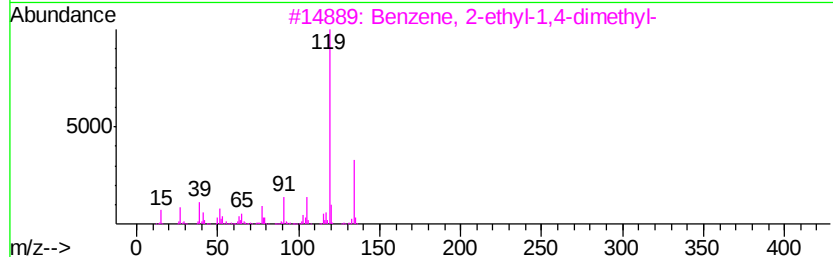
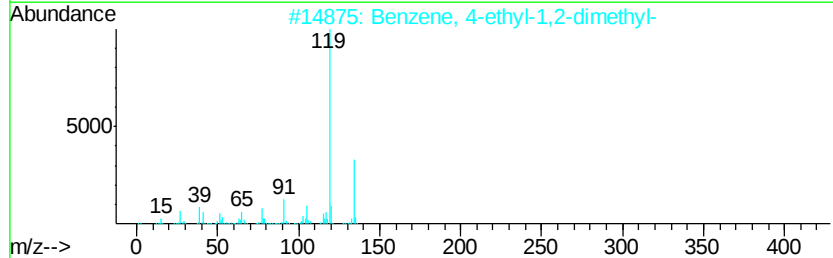
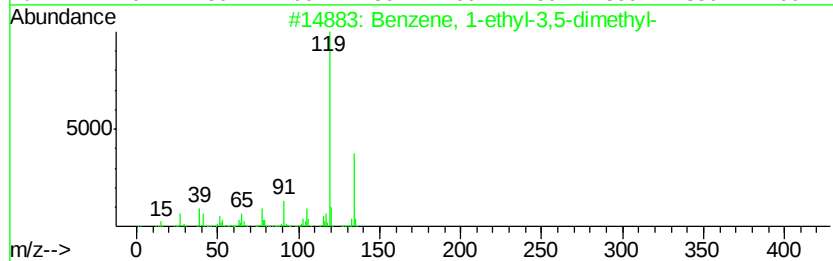
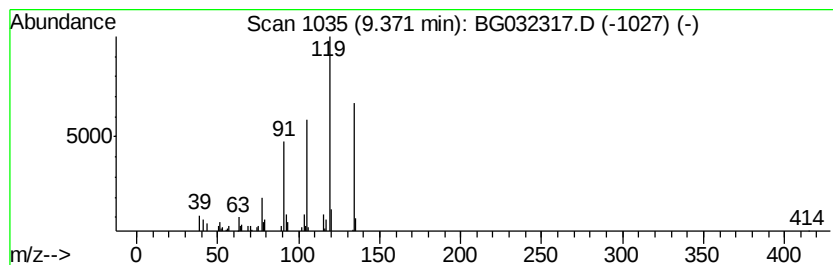
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 Benzene, 1-ethyl-3,5-dimethyl- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.37	5.77 ng	52646	1,4-Dichlorobenzene-d4	8.72

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	94
2		Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	91
3		Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	91
4		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	87
5		Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	87



Data Path : Z:\HPCHEM1\BNA G\DATA\BG012518\
Data File : BG032317.D
Acq On : 26 Jan 2018 6:39
Operator : SJ/JU
Sample : J1267-03
Misc :
ALS Vial : 22 Sample Multiplier: 1

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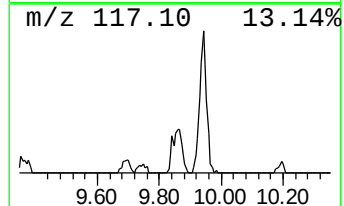
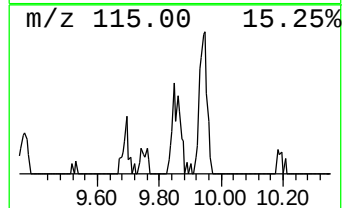
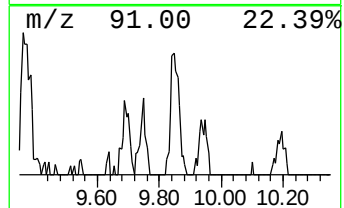
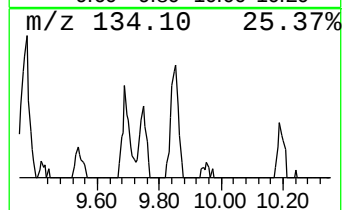
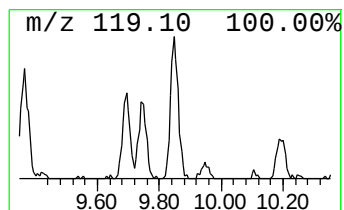
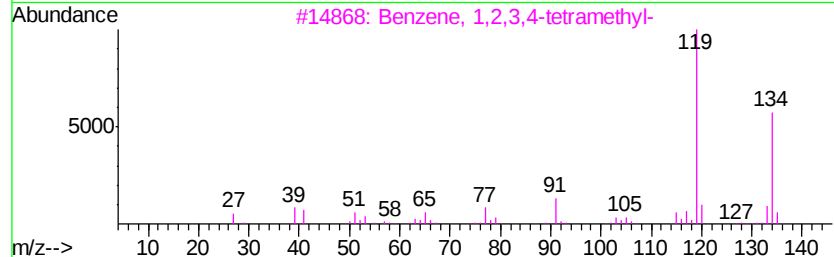
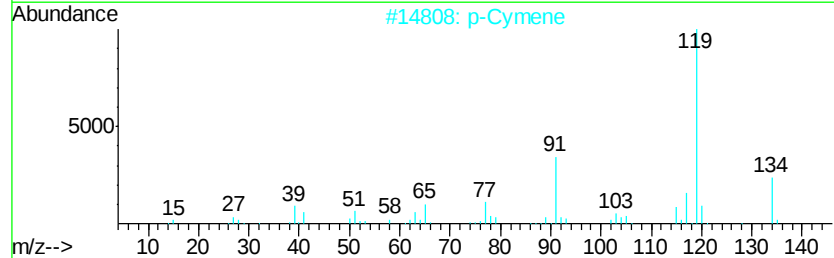
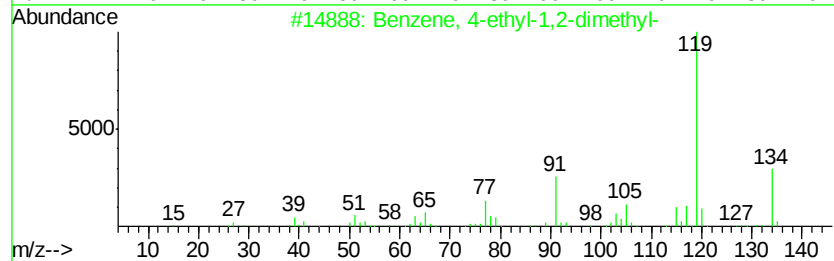
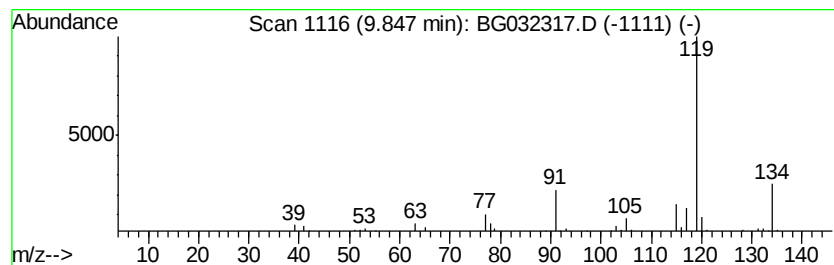
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.85	5.56 ng	50771	1,4-Dichlorobenzene-d4	8.72

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	87
2			p-Cymene	134	C10H14	000099-87-6	83
3			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	80
4			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	80
5			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	80



Data Path : Z:\HPCHEM1\BNA G\DATA\BG012518\
Data File : BG032317.D
Acq On : 26 Jan 2018 6:39
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Sample : J1267-03
Misc :
ALS Vial : 22 Sample Multiplier: 1

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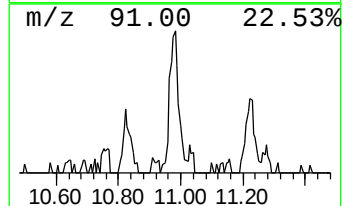
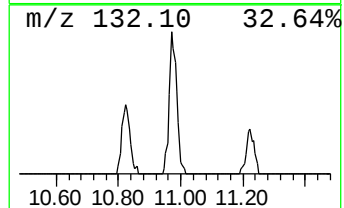
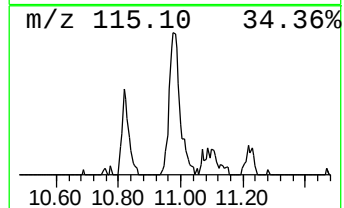
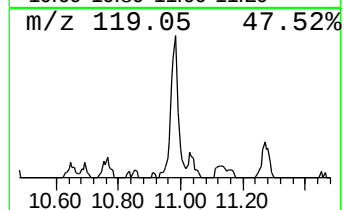
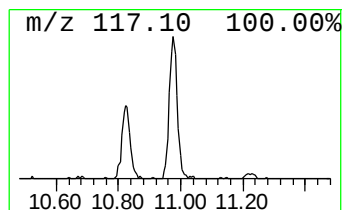
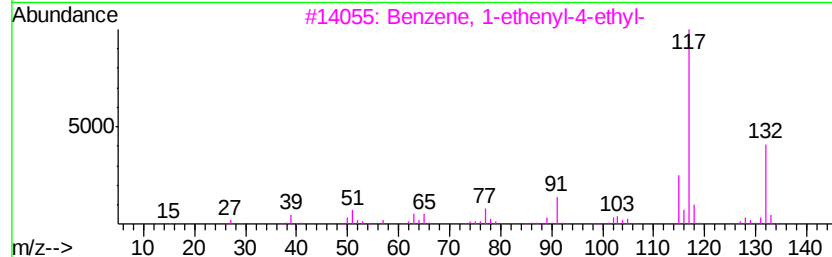
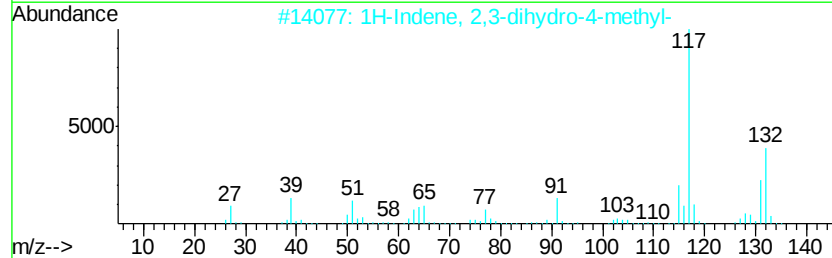
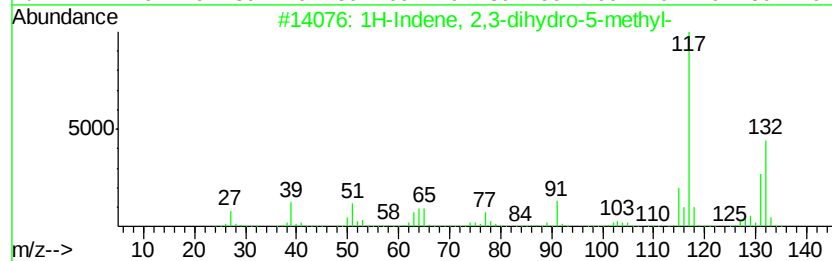
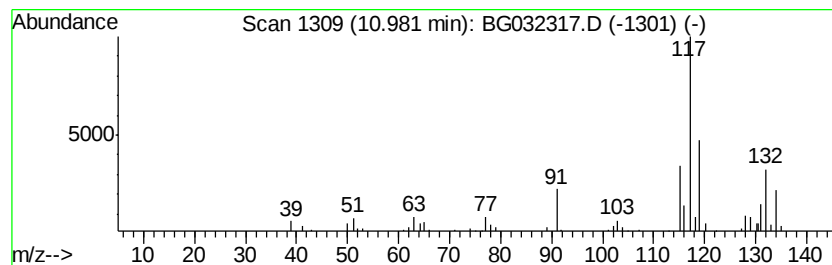
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 1H-Indene, 2,3-dihydro-5-me... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.98	9.00 ng	112632	Napthalene-d8	11.60

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	83
2		1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	83
3		Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	76
4		Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	74
5		3-Phenylbut-1-ene	132	C10H12	000934-10-1	74



Data Path : Z:\HPCHEM1\BNA G\DATA\BG012518\
Data File : BG032317.D
Acq On : 26 Jan 2018 6:39
Operator : SJ/JU
Sample : J1267-03
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
IGW-6

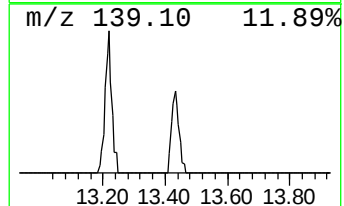
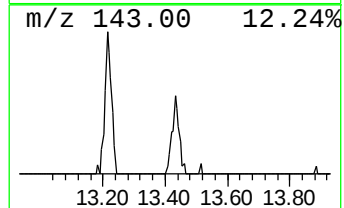
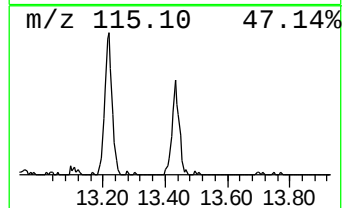
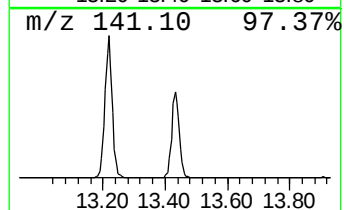
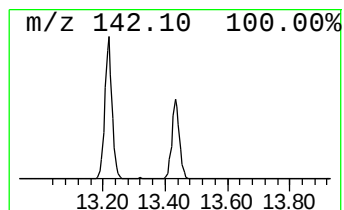
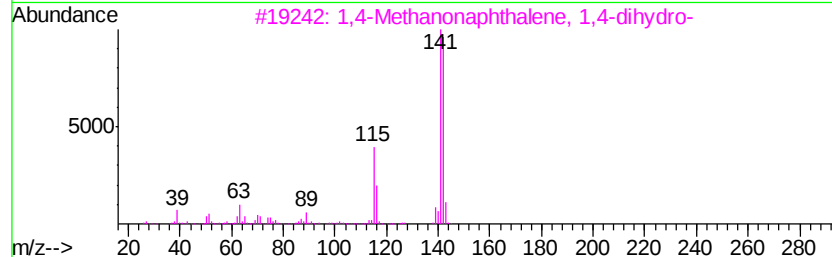
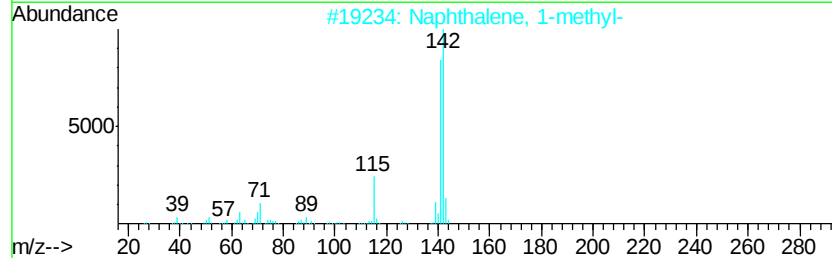
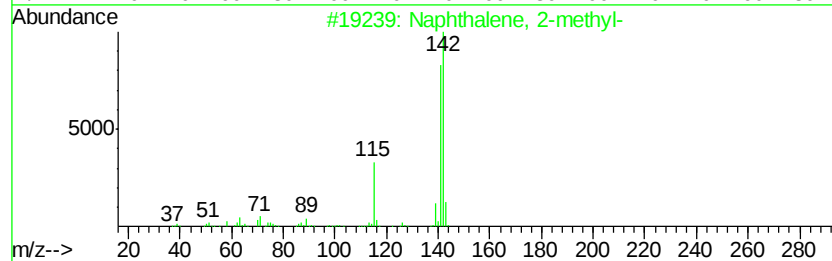
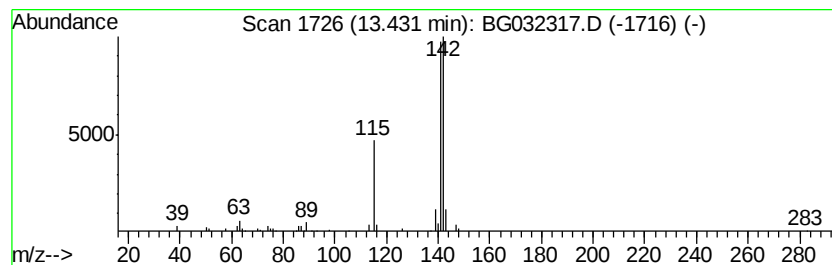
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 19 Naphthalene, 1-methyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.43	7.64 ng	95629	Naphthalene-d8	11.60

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2-methyl-	142	C11H10	000091-57-6	94
2			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	94
3			1,4-Methanonaphthalene, 1,4-dihv...	142	C11H10	004453-90-1	91
4			Benzocycloheptatriene	142	C11H10	000264-09-5	91
5			Spiro(9-methylenetricyclo[6.2.1....	184	C13H12O	033929-47-4	72



Data Path : Z:\HPCHEM1\BNA G\DATA\BG012518\
Data File : BG032317.D
Acq On : 26 Jan 2018 6:39
Operator : SJ/JU
Sample : J1267-03
Misc :
ALS Vial : 22 Sample Multiplier: 1

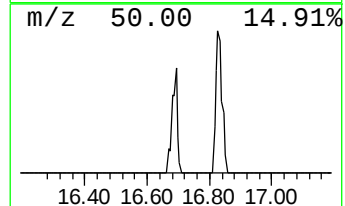
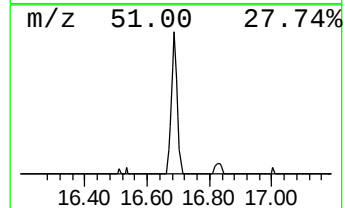
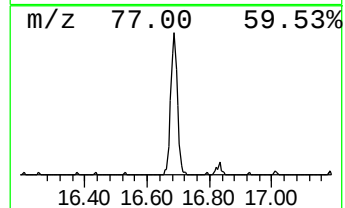
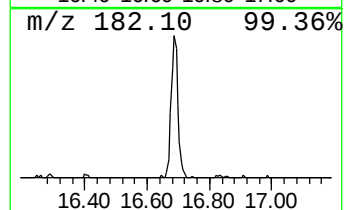
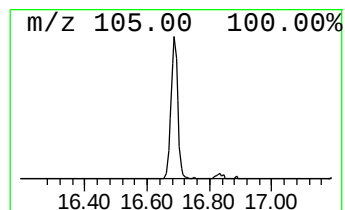
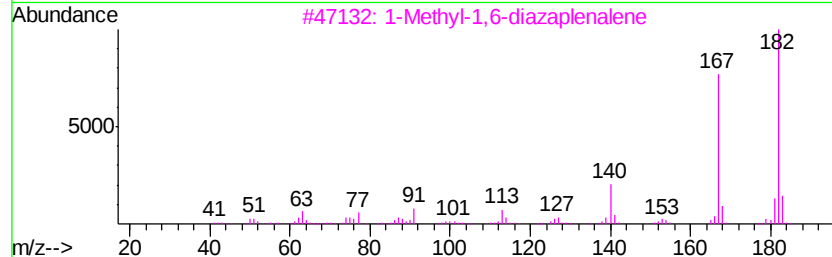
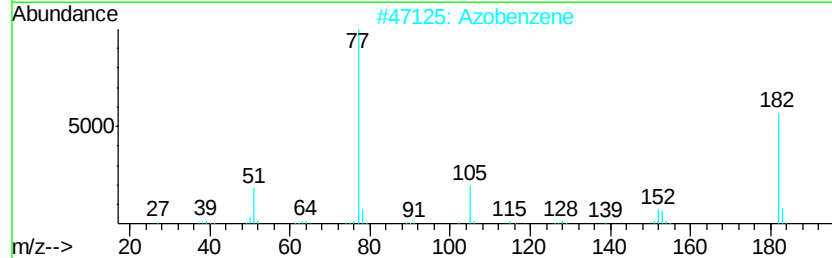
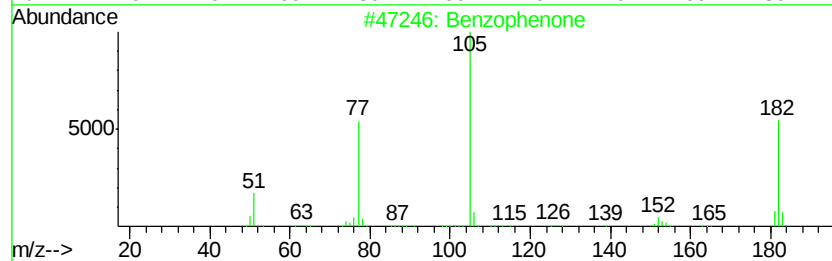
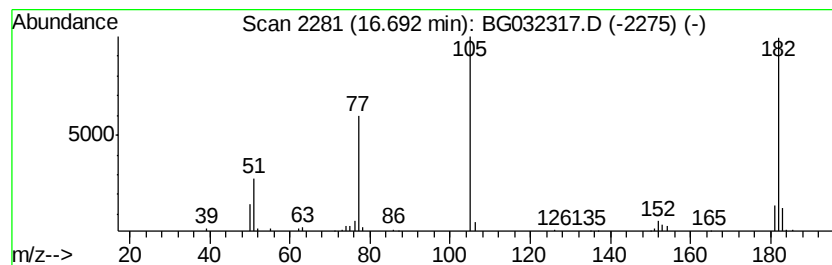
Instrument :
BNA_G
ClientSampleId :
IGW-6

Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG011218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 20 Benzophenone Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.	
16.69	4.06 ng	86379	Acenaphthene-d10	15.36	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzophenone	182	C13H10O	000119-61-9	97
2	Azobenzene	182	C12H10N2	000103-33-3	50
3	1-Methyl-1,6-diazaplenalene	182	C12H10N2	079691-99-9	43
4	Benzaldehyde, p-chloro-, dimethy...	182	C9H11ClN2	022699-29-2	43
5	2,4,6-Trimethoxytoluene	182	C10H14O3	014107-97-2	43



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG012518\
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 Sample : J1267-03
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 IGW-6

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG011218.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
Butane, 2-methoxy...	3.47	46.3	ng	422600	1	8.72	182474	20.0
2-Pentanone, 4-hy...	5.65	97.1	ng	885692	1	8.72	182474	20.0
Benzene, 1,3-dime...	6.23	53.0	ng	483896	1	8.72	182474	20.0
Benzene, propyl-	7.65	4.0	ng	36744	1	8.72	182474	20.0
Benzene, 1-ethyl-...	7.76	15.2	ng	138590	1	8.72	182474	20.0
Benzene, 1,2,4-tr...	7.90	10.4	ng	94880	1	8.72	182474	20.0
unknown8.24	8.24	74.2	ng	677182	1	8.72	182474	20.0
Benzene, 1,2,3-tr...	8.35	39.9	ng	363536	1	8.72	182474	20.0
Benzene, 1,1'-(1-...	9.11	10.8	ng	98884	1	8.72	182474	20.0
Benzene, 1-ethyl-...	9.37	5.8	ng	52646	1	8.72	182474	20.0
Benzene, 4-ethyl-...	9.85	5.6	ng	50771	1	8.72	182474	20.0
1H-Indene, 2,3-di...	10.98	9.0	ng	112632	2	11.60	250390	20.0
Naphthalene, 1-me...	13.43	7.6	ng	95629	2	11.60	250390	20.0
Benzophenone	16.69	4.1	ng	86379	3	15.36	425379	20.0