

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA G\DATA\BG111317\
 Data File : BG030984.D
 Acq On : 13 Nov 2017 22:02
 Operator : SJ/JU
 Sample : I6436-03
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 153RD-63RD-ST-DRUM

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-BG111017.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	5.220	322	329	341	rBV	73267	118074	1.38%	0.321%
2	5.519	374	380	391	rBV	104714	155070	1.82%	0.422%
3	5.631	393	399	406	rVV	339719	551132	6.46%	1.500%
4	5.707	406	412	416	rVV	380360	609745	7.14%	1.659%
5	5.772	416	423	439	rVB	3613977	8537924	100.00%	23.230%
6	6.142	477	486	494	rBV	2610024	4354549	51.00%	11.848%
7	7.223	660	670	675	rBV	168200	276677	3.24%	0.753%
8	7.305	675	684	689	rVV2	230642	525816	6.16%	1.431%
9	7.358	689	693	699	rVB	104017	172080	2.02%	0.468%
10	7.523	715	721	732	rVB	107521	177621	2.08%	0.483%
11	7.676	740	747	757	rBV	711142	1268832	14.86%	3.452%
12	7.787	759	766	773	rVB	282154	478788	5.61%	1.303%
13	8.140	819	826	840	rBV	160213	305153	3.57%	0.830%
14	8.257	840	846	854	rVB	116779	200793	2.35%	0.546%
15	8.469	875	882	888	rBV	729321	1280784	15.00%	3.485%
16	8.780	929	935	939	rVV	50634	87047	1.02%	0.237%
17	9.309	1018	1025	1035	rVB	548217	988461	11.58%	2.689%
18	9.785	1095	1106	1111	rBV2	143687	384809	4.51%	1.047%
19	10.954	1298	1305	1313	rBV	223253	419123	4.91%	1.140%
20	11.477	1385	1394	1400	rBV2	113928	266551	3.12%	0.725%
21	11.759	1435	1442	1456	rBV3	52501	151794	1.78%	0.413%
22	11.888	1457	1464	1476	rVB5	36836	112852	1.32%	0.307%
23	12.564	1572	1579	1587	rBV	106749	205462	2.41%	0.559%
24	13.381	1708	1718	1728	rBV	1759497	2805269	32.86%	7.633%
25	14.197	1849	1857	1867	rBV	171374	283204	3.32%	0.771%
26	14.756	1945	1952	1961	rVB2	389472	663636	7.77%	1.806%
27	16.089	2172	2179	2187	rBV2	247564	373007	4.37%	1.015%
28	16.242	2198	2205	2213	rBV	1590882	2479847	29.05%	6.747%
29	16.342	2218	2222	2225	rVV2	68883	109856	1.29%	0.299%
30	16.383	2225	2229	2234	rVV3	57738	116207	1.36%	0.316%
31	16.571	2254	2261	2270	rBV2	79828	191263	2.24%	0.520%
32	17.494	2406	2418	2424	rBV2	574456	900019	10.54%	2.449%
33	17.781	2461	2467	2474	rBV	124467	208395	2.44%	0.567%
34	20.085	2853	2859	2865	rBV	3441802	4785821	56.05%	13.021%

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Instrument :
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ClientSampleId :
153RD-63RD-ST-DRUM

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-BG111017.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	21.759	3137	3144	3157	rBV2	657326	1119209	13.11%	3.045%
36	25.055	3693	3705	3716	rBV2	360474	1088409	12.75%	2.961%

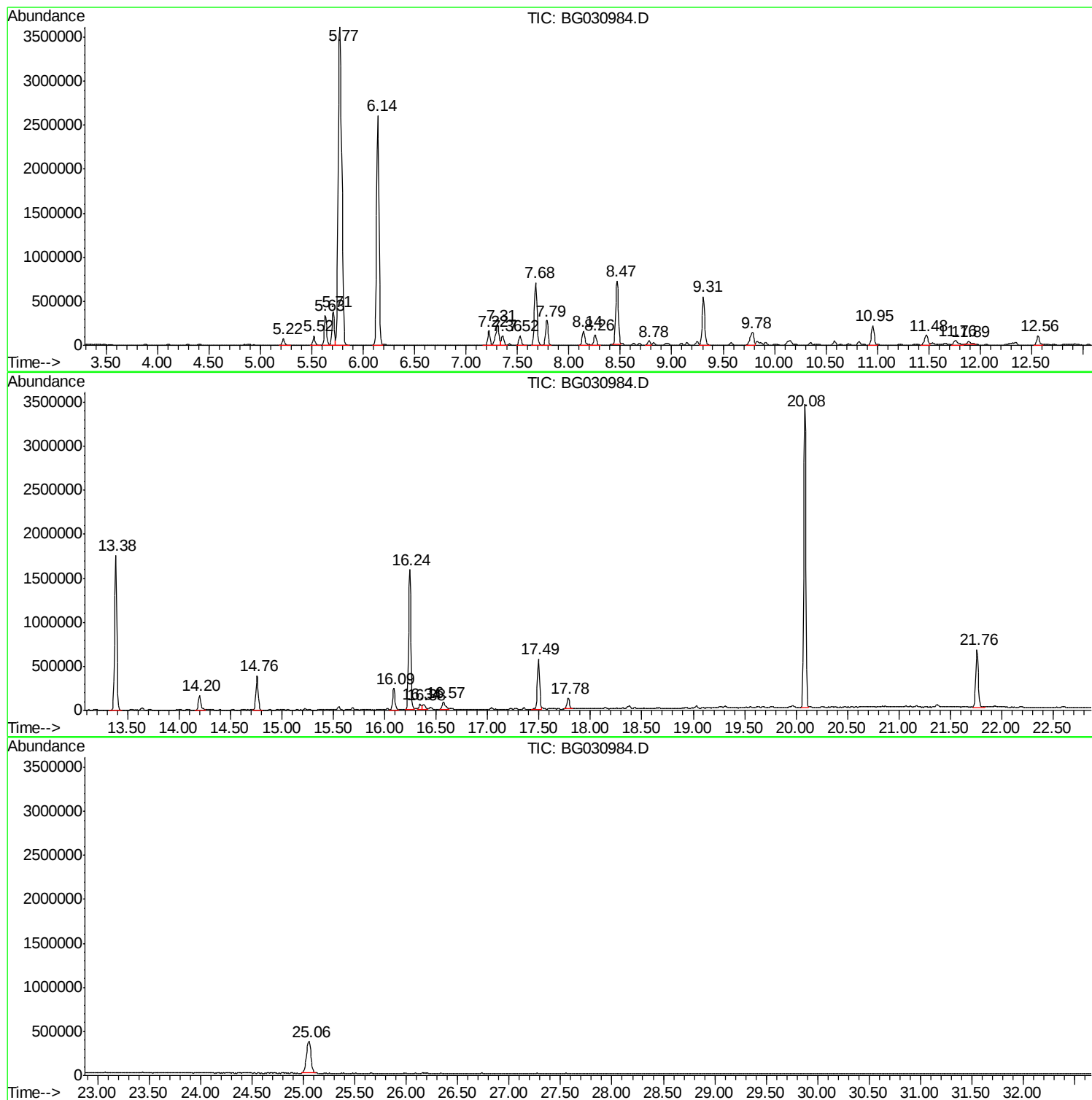
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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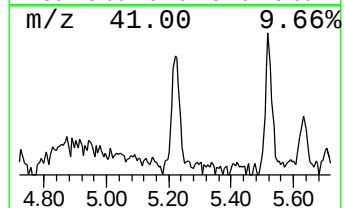
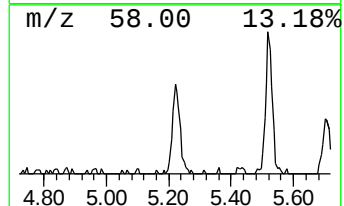
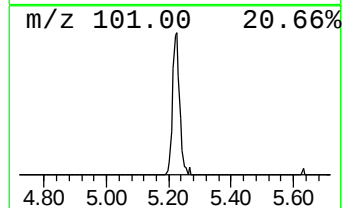
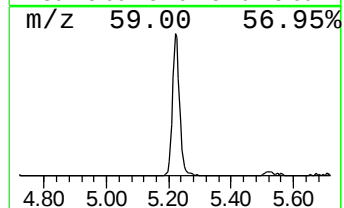
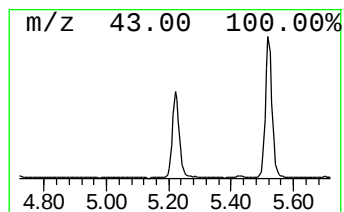
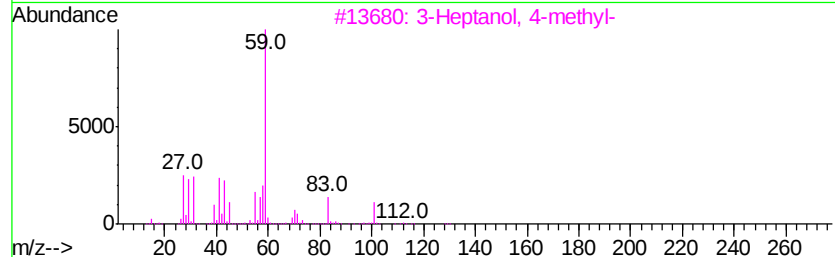
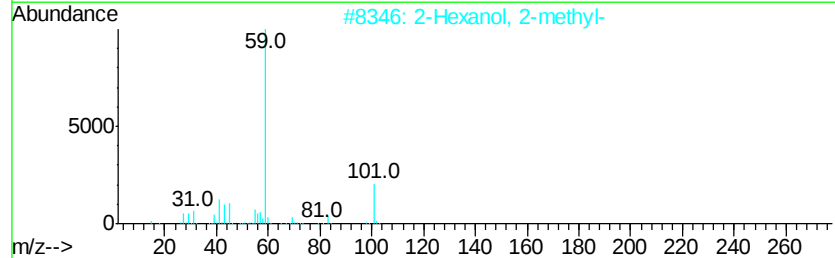
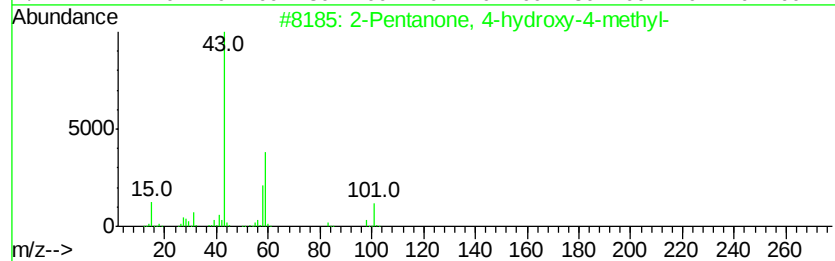
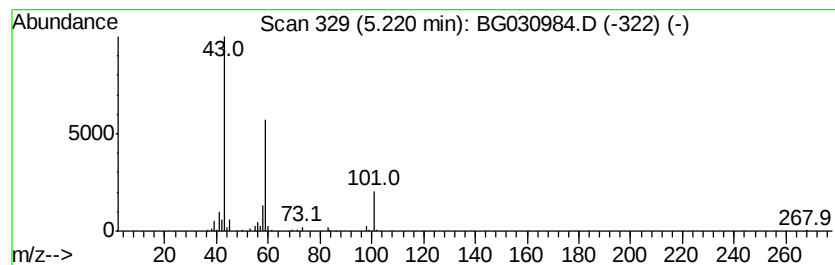
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.22	7.74 ng	118074	1,4-Dichlorobenzene-d4	8.14

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	39
2			2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	28
3			3-Heptanol, 4-methyl-	130	C8H18O	014979-39-6	25
4			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	25
5			Allyl dithioacetate	132	C5H8S2	027249-83-8	23



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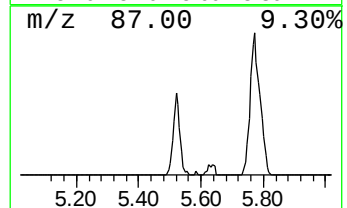
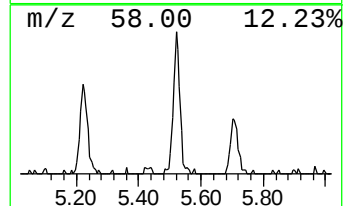
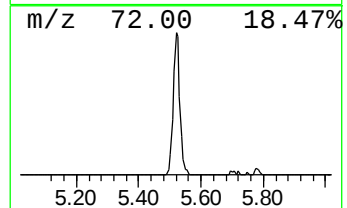
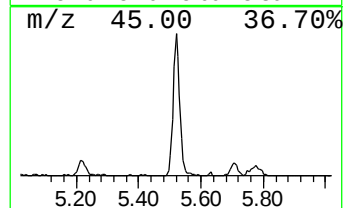
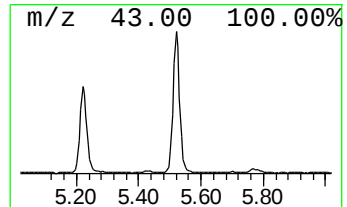
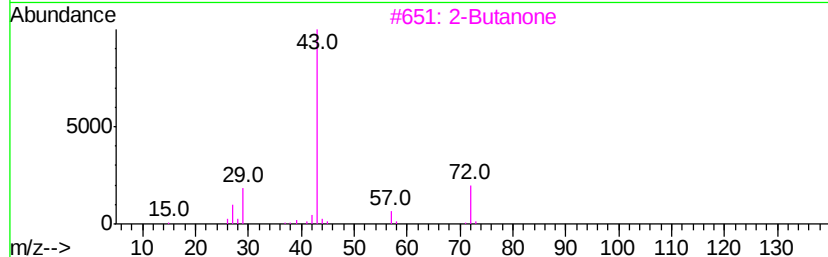
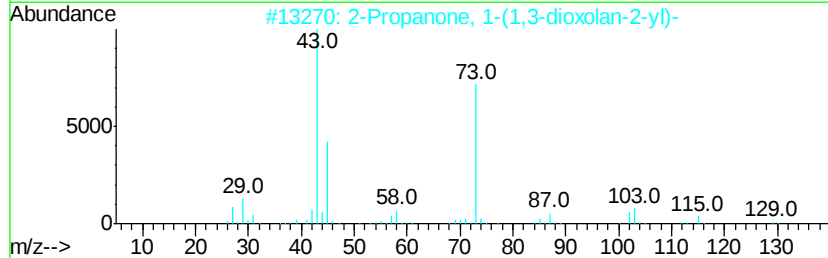
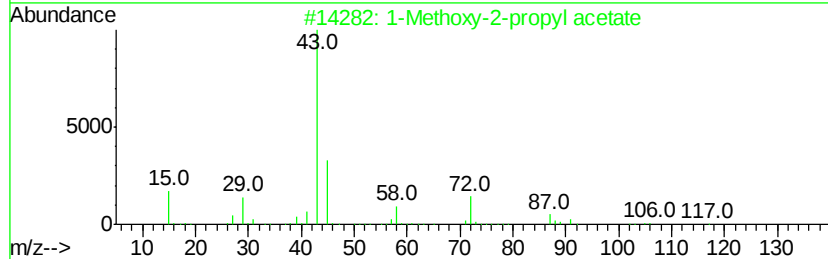
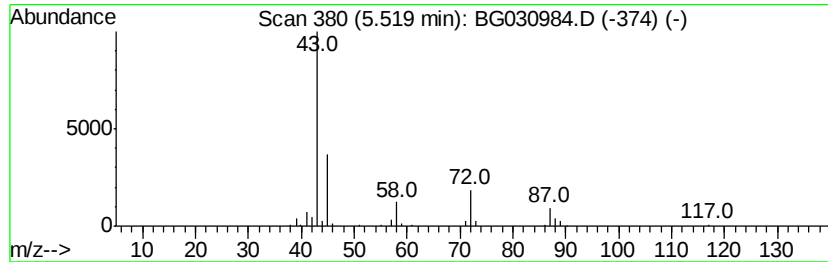
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 1-Methoxy-2-propyl acetate Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.52	10.16 ng	155070	1,4-Dichlorobenzene-d4	8.14

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Methoxy-2-propyl acetate	132	C6H12O3	000108-65-6	78
2		2-Propanone, 1-(1,3-dioxolan-2-yl)-	130	C6H10O3	000767-04-4	36
3		2-Butanone	72	C4H8O	000078-93-3	25
4		1,2-Propanediol, 2-acetate	118	C5H10O3	006214-01-3	23
5		Propane, 1-(1-methylethoxy)-	102	C6H14O	000627-08-7	23



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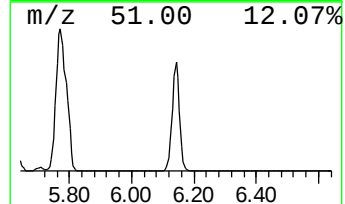
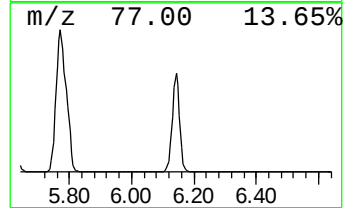
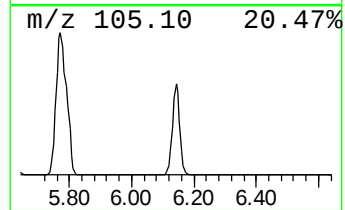
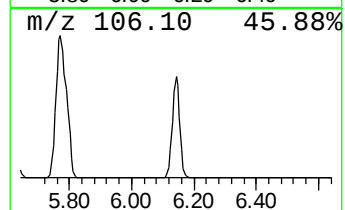
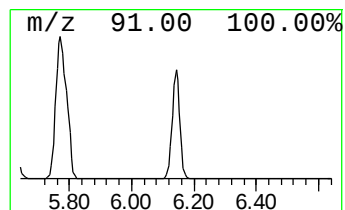
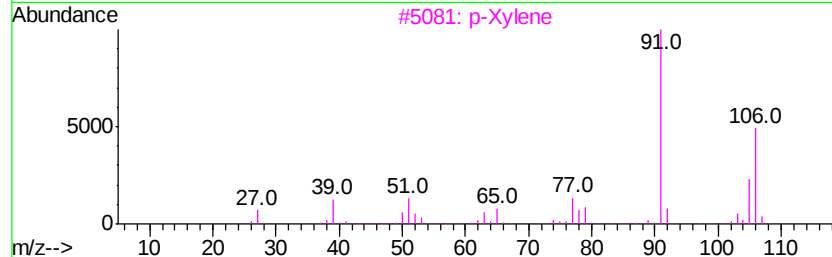
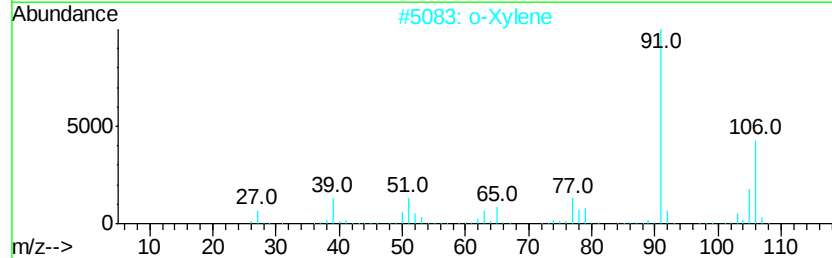
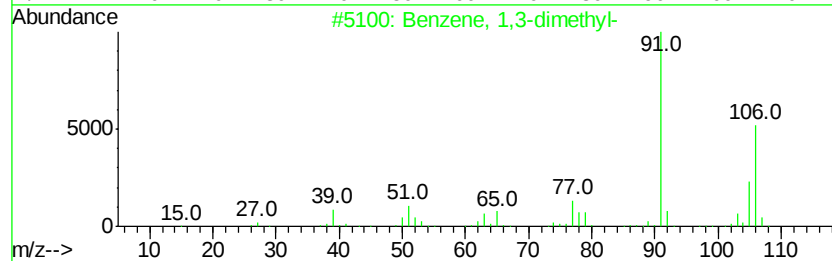
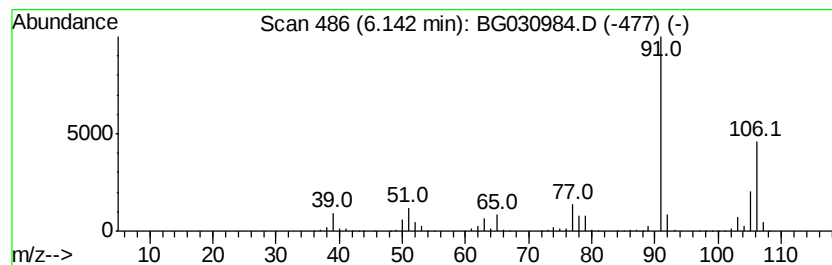
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Benzene, 1,3-dimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.14	285.40 ng	4354550	1,4-Dichlorobenzene-d4	8.14

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	87
5		Cyclopentene, 1-ethenyl-3-methyl...	106	C8H10	061142-07-2	86



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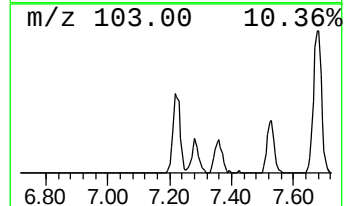
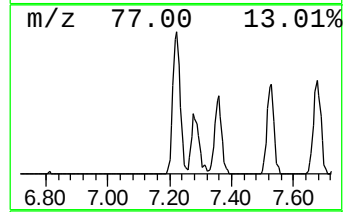
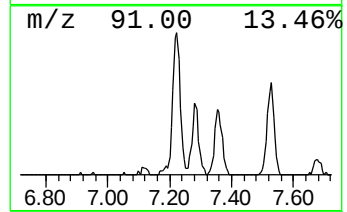
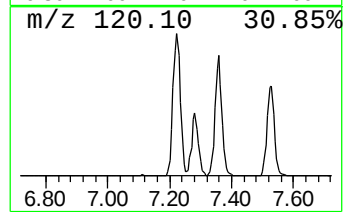
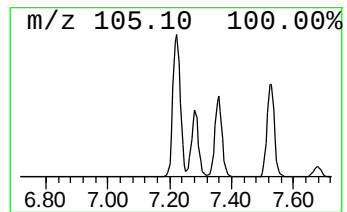
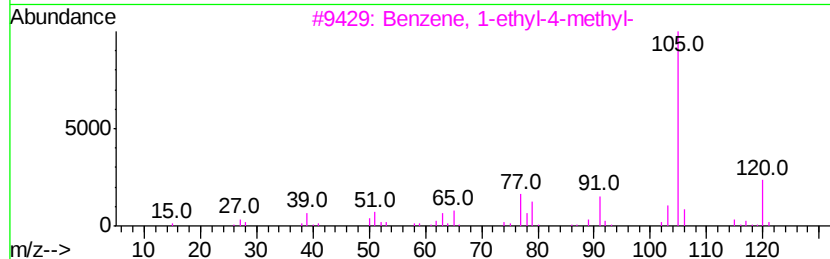
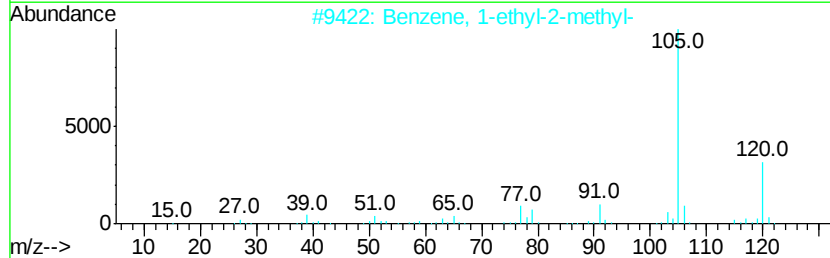
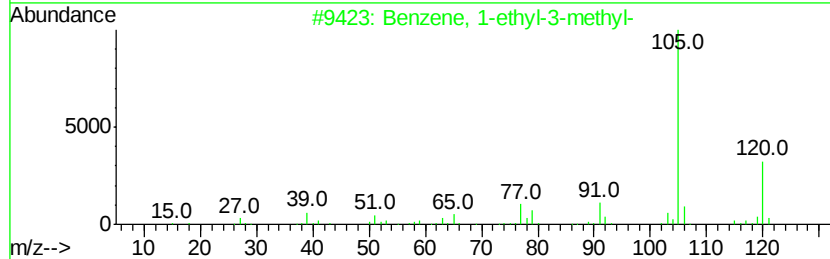
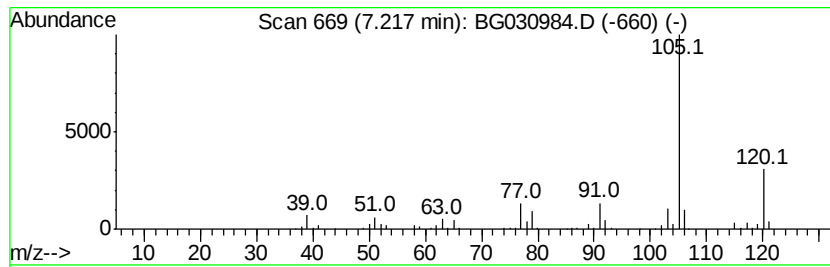
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Benzene, 1-ethyl-3-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.22	18.13 ng	276677	1,4-Dichlorobenzene-d4	8.14

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	97
2		Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
3		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	93
4		Mesitylene	120	C9H12	000108-67-8	91
5		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91



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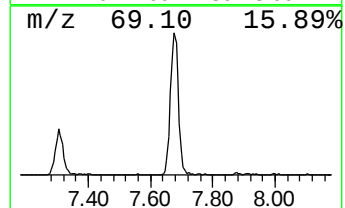
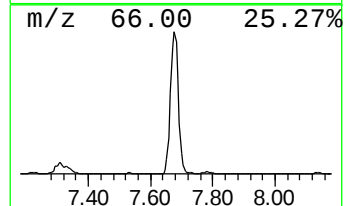
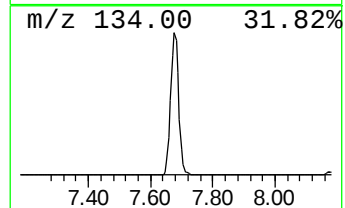
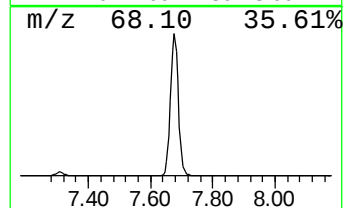
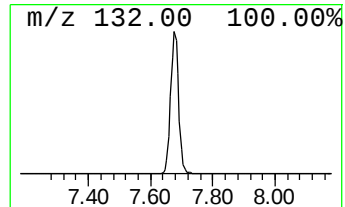
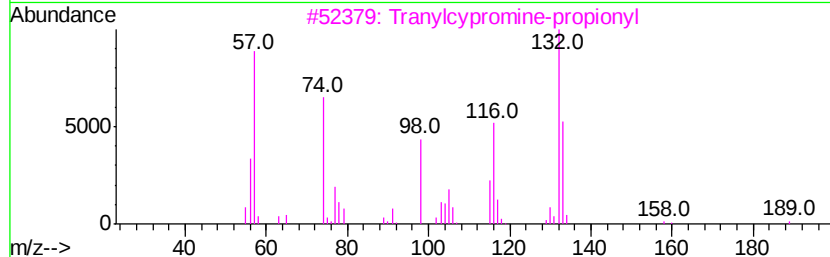
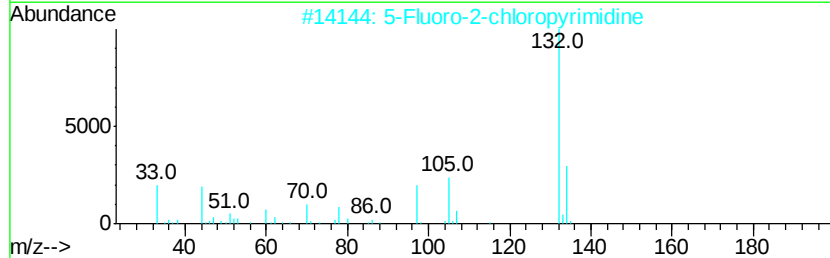
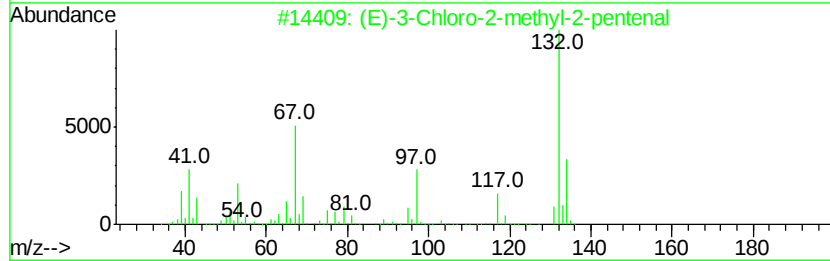
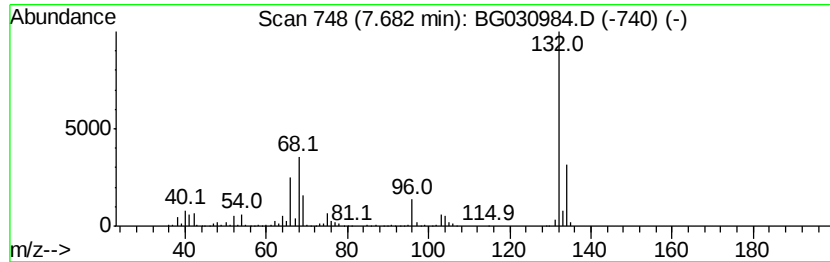
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TIC Library : C:\Database\NIST11.L
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Peak Number 8 unknown7.68 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.68	83.16 ng	1268830	1,4-Dichlorobenzene-d4	8.14

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	10
4		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
5		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9



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Acq On : 13 Nov 2017 22:02
Operator : SJ/JU
Sample : I6436-03
Misc :
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Instrument :
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ClientSampleId :
153RD-63RD-ST-DRUM

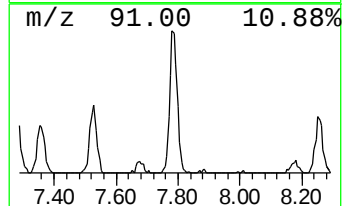
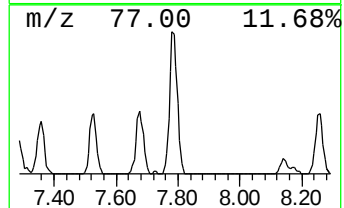
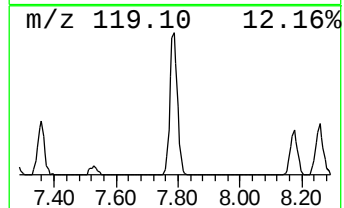
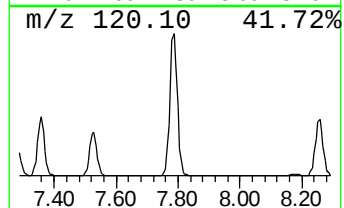
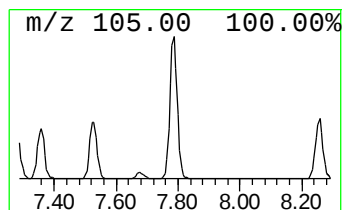
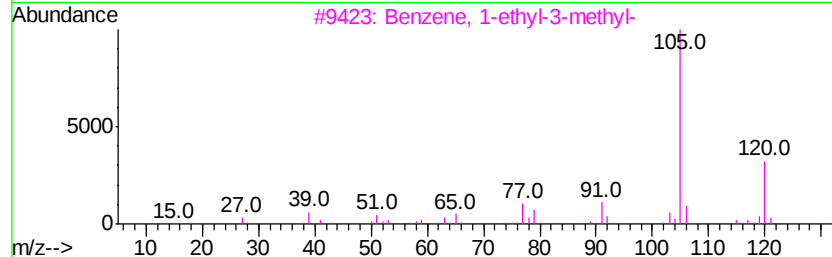
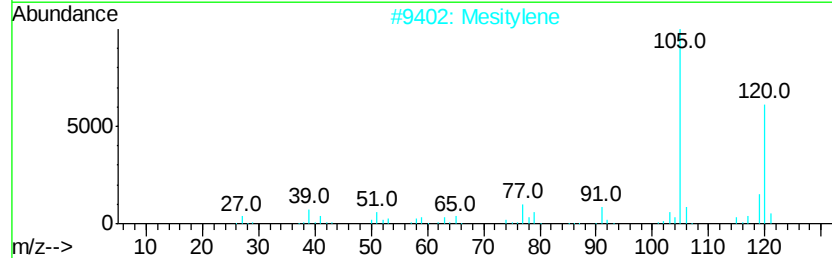
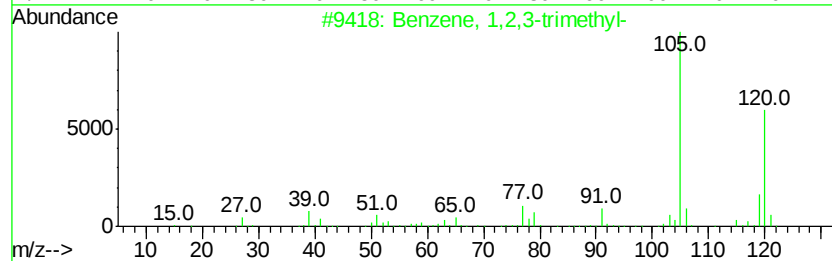
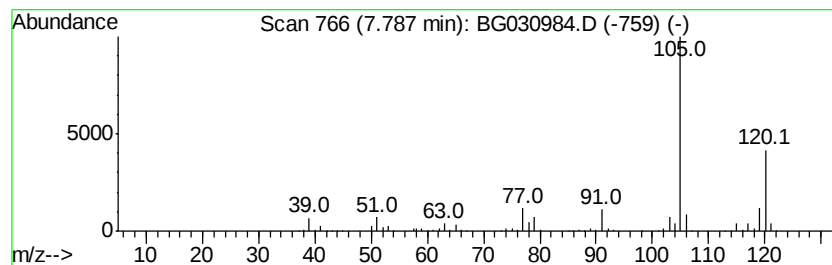
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Benzene, 1,2,3-trimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.79	31.38 ng	478788	1,4-Dichlorobenzene-d4	8.14

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	97
2		Mesitylene	120	C9H12	000108-67-8	97
3		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
4		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
5		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91



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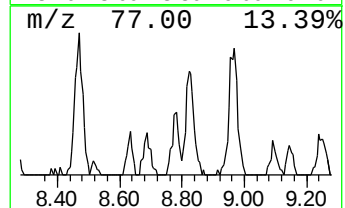
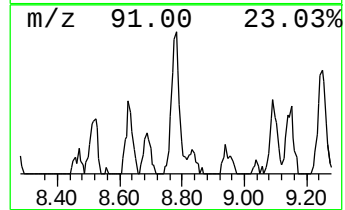
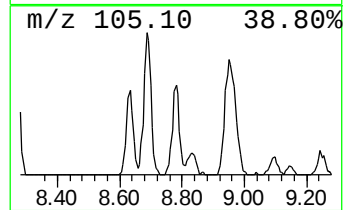
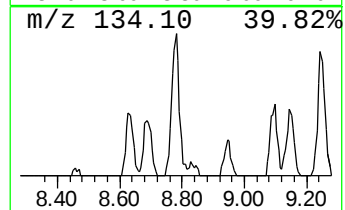
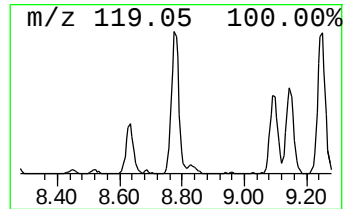
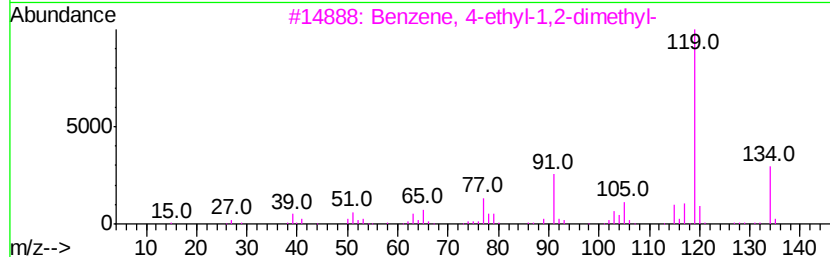
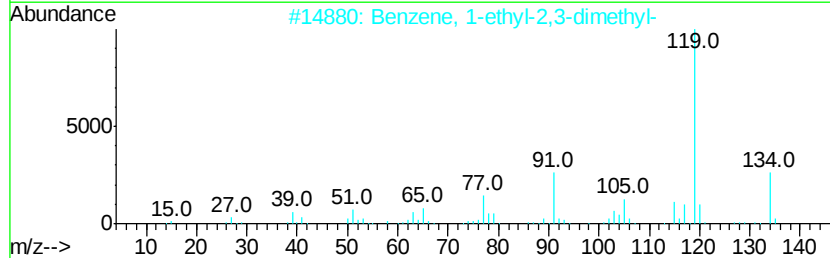
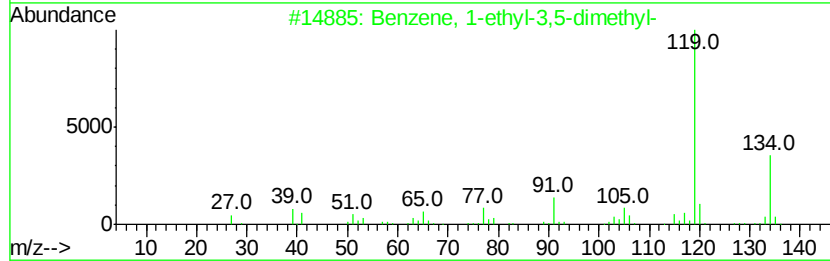
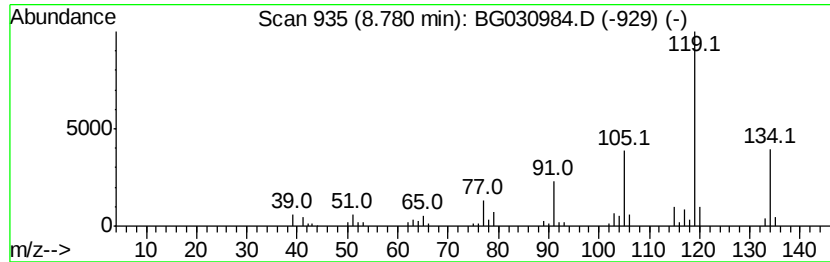
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Benzene, 1-ethyl-3,5-dimethyl- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.78	5.71 ng	87047	1,4-Dichlorobenzene-d4	8.14

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	95
2		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	93
3		Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	93
4		Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	93
5		Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	91



Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA G\DATA\BG111317\
Data File : BG030984.D
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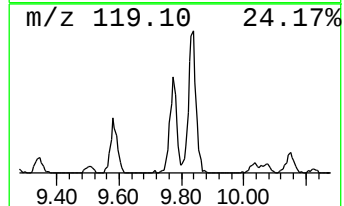
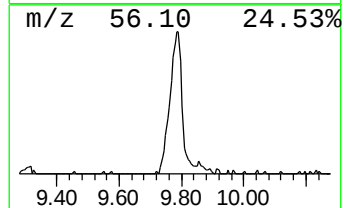
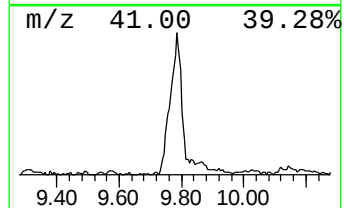
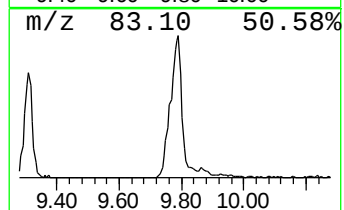
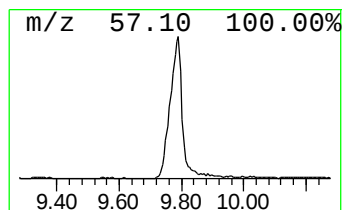
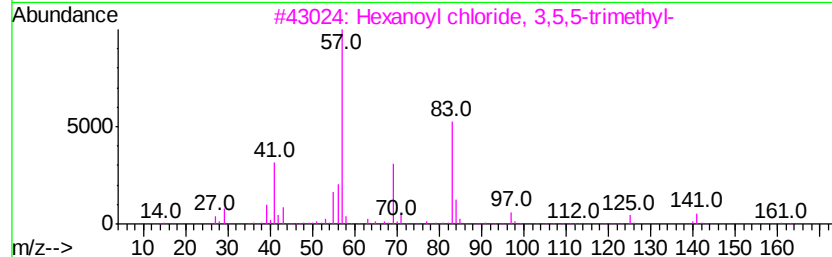
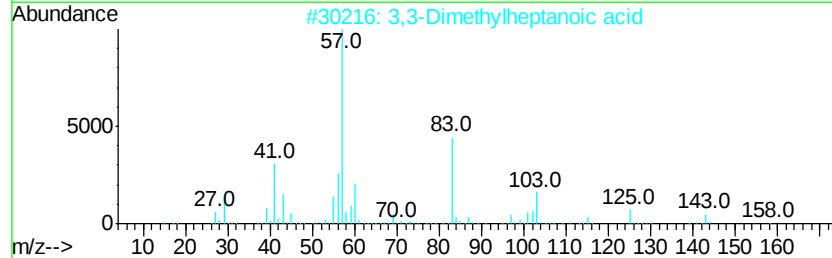
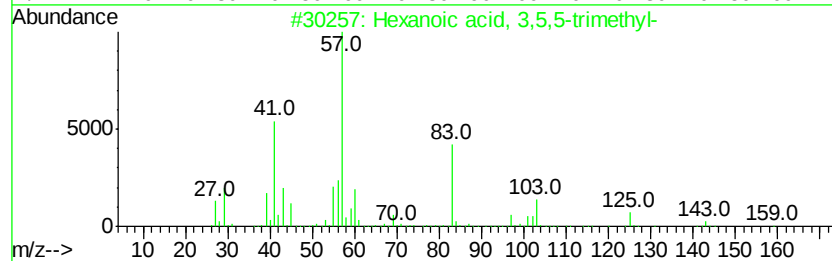
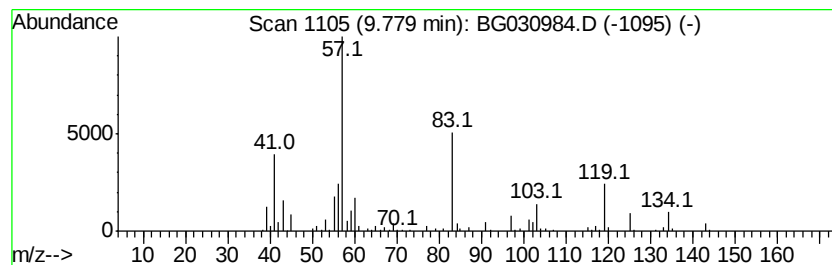
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Hexanoic acid, 3,5,5-trimet... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.78	18.36 ng	384809	Napthalene-d8	10.95

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexanoic acid, 3,5,5-trimethyl-	158	C9H18O2	003302-10-1	80
2		3,3-Dimethylheptanoic acid	158	C9H18O2	067061-30-7	72
3		Hexanoyl chloride, 3,5,5-trimethyl-	176	C9H17ClO	036727-29-4	50
4		4,4-Dimethyl-2-pentanol, chlorod...	228	C9H15ClF2O2	1000376-27-0	38
5		Hexanal, 3,5,5-trimethyl-	142	C9H18O	005435-64-3	38



Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA G\DATA\BG111317\
Data File : BG030984.D
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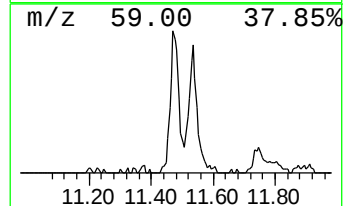
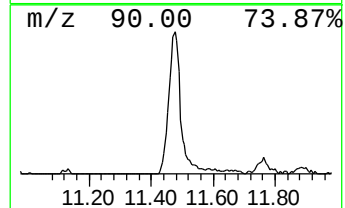
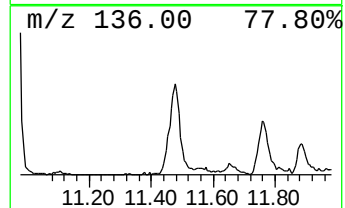
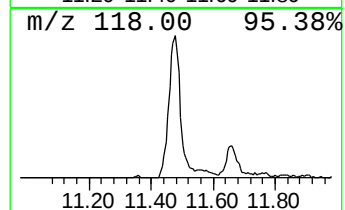
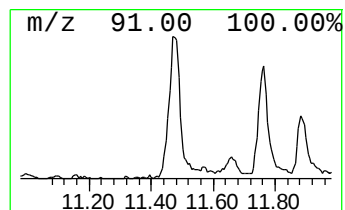
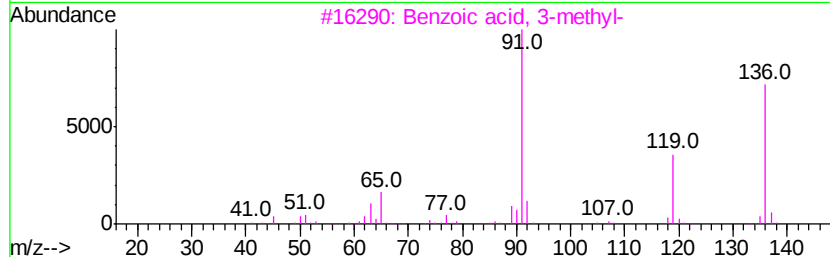
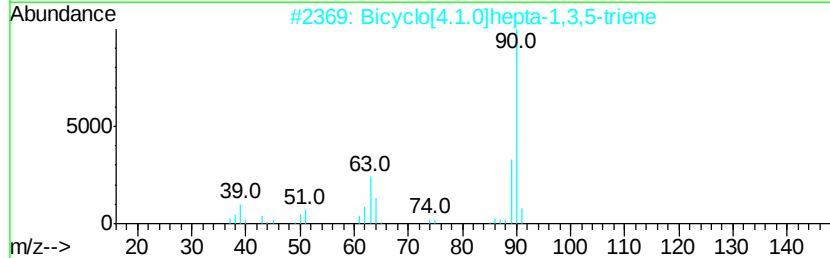
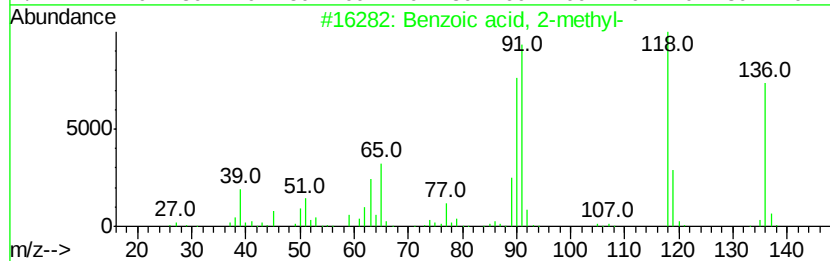
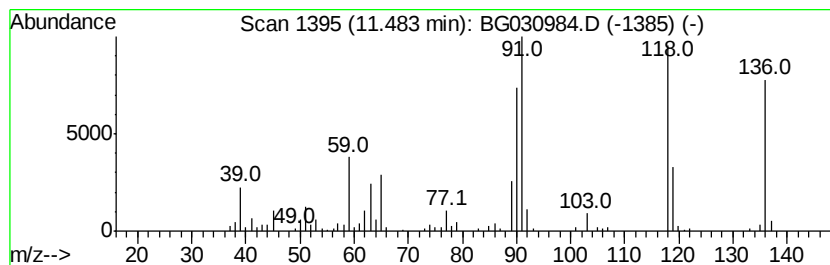
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Benzoic acid, 2-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.48	12.72 ng	266551	Naphthalene-d8	10.95

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzoic acid, 2-methyl-	136	C8H8O2	000118-90-1	98
2		Bicyclo[4.1.0]hepta-1,3,5-triene	90	C7H6	004646-69-9	43
3		Benzoic acid, 3-methyl-	136	C8H8O2	000099-04-7	35
4		Benzoic acid, 4-methyl-	136	C8H8O2	000099-94-5	25
5		1H-Indole, 2,3-dihydro-	119	C8H9N	000496-15-1	15



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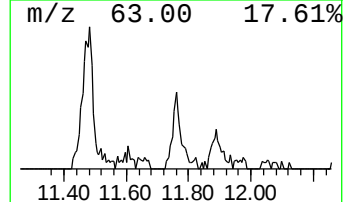
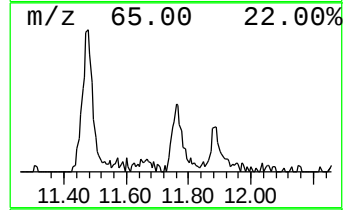
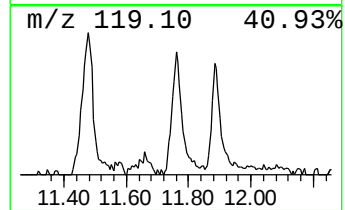
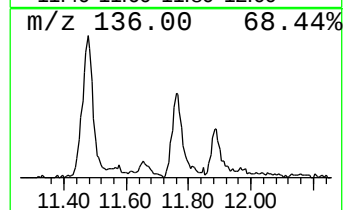
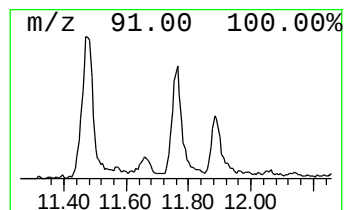
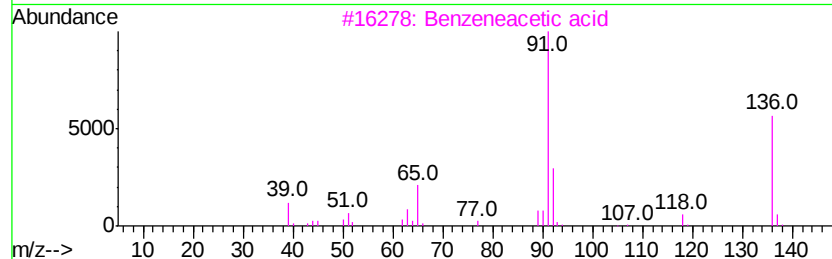
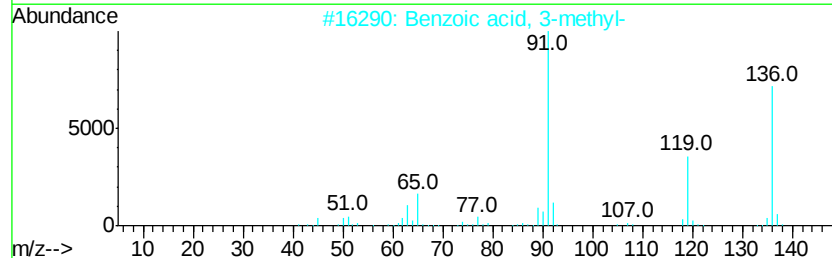
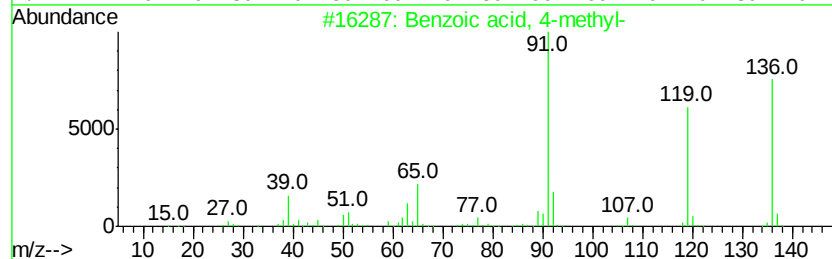
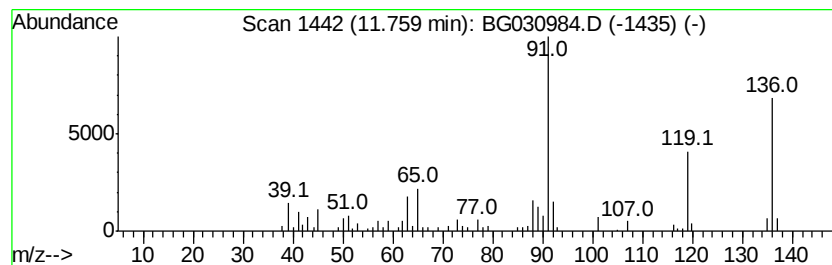
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Benzoic acid, 4-methyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.76	7.24 ng	151794	Naphthalene-d8	10.95

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzoic acid, 4-methyl-	136	C8H8O2	000099-94-5	93
2		Benzoic acid, 3-methyl-	136	C8H8O2	000099-04-7	93
3		Benzeneacetic acid	136	C8H8O2	000103-82-2	55
4		Ethanone, 1-(3-methylphenyl)-	134	C9H10O	000585-74-0	47
5		m-Toluamide	135	C8H9NO	000618-47-3	43



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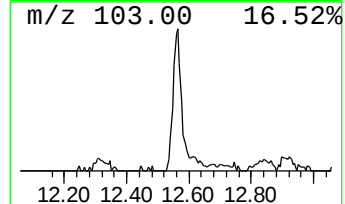
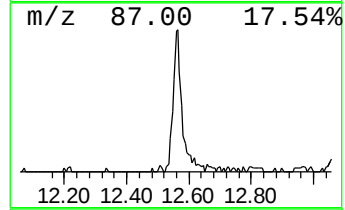
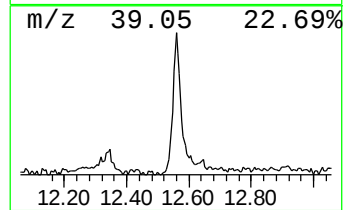
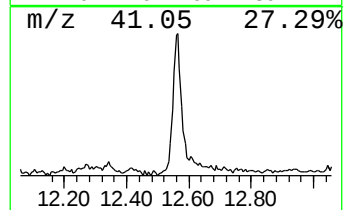
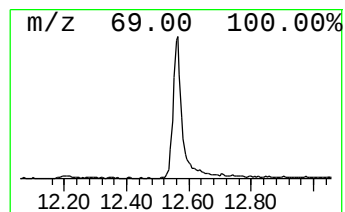
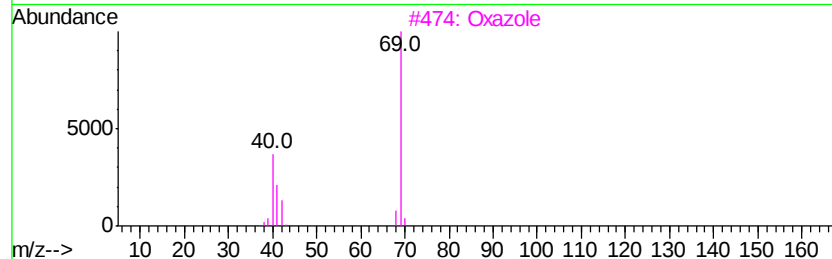
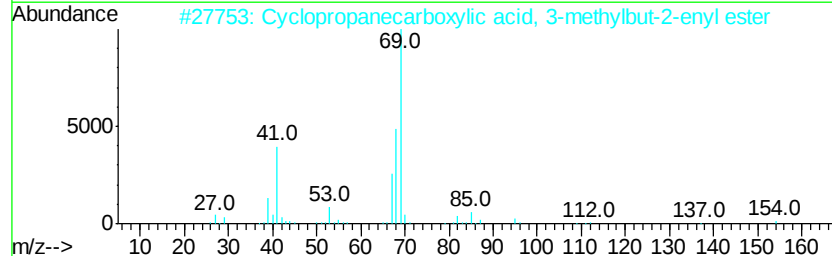
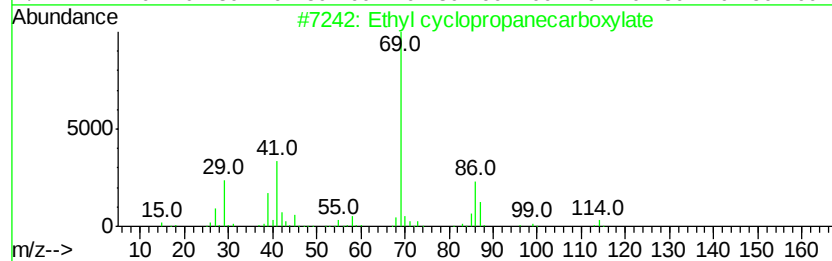
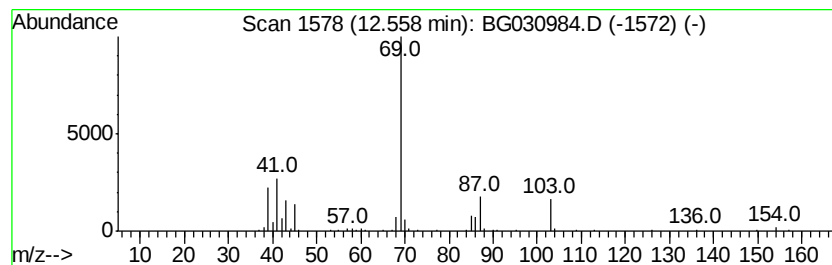
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 unknown12.56 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.56	9.80 ng	205462	Naphthalene-d8	10.95

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethyl cyclopropanecarboxylate	114	C6H10O2	004606-07-9	39
2			Cyclopropanecarboxylic acid, 3-m...	154	C9H14O2	1000299-37-4	9
3			Oxazole	69	C3H3NO	000288-42-6	9
4			Crotonic anhydride	154	C8H10O3	000623-68-7	9
5			2-Butenoic acid, methyl ester	100	C5H8O2	018707-60-3	9



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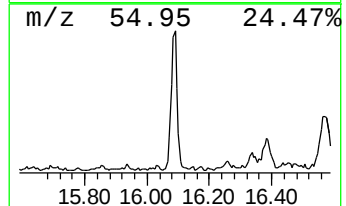
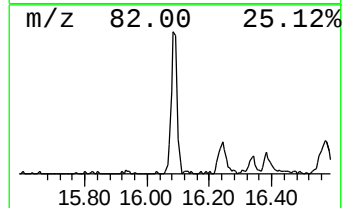
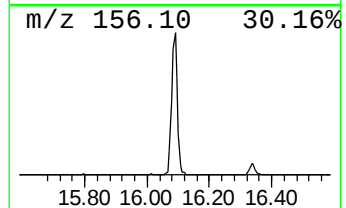
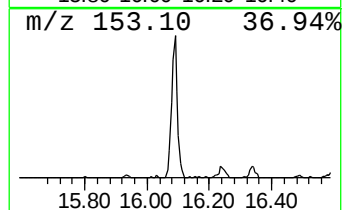
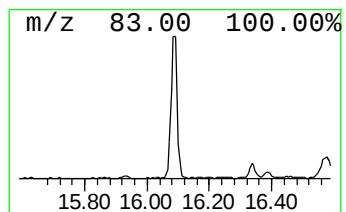
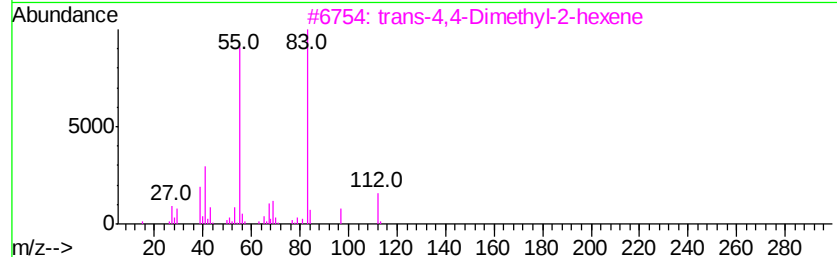
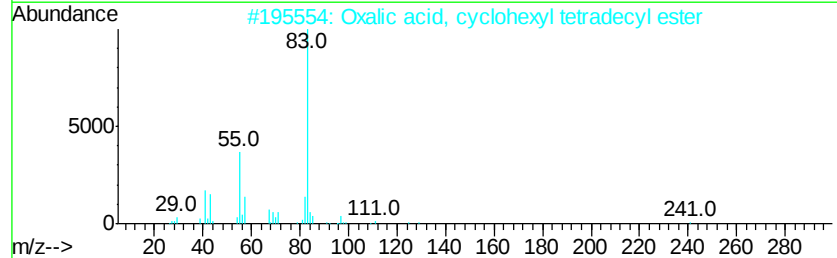
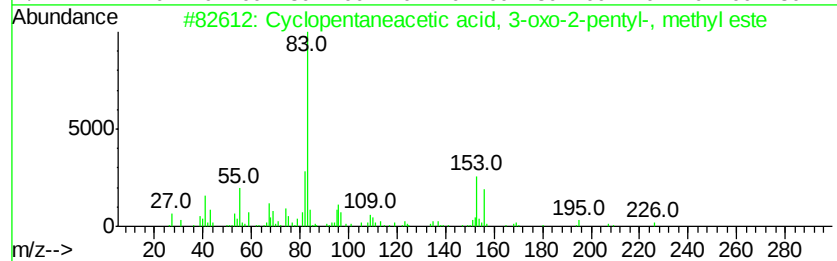
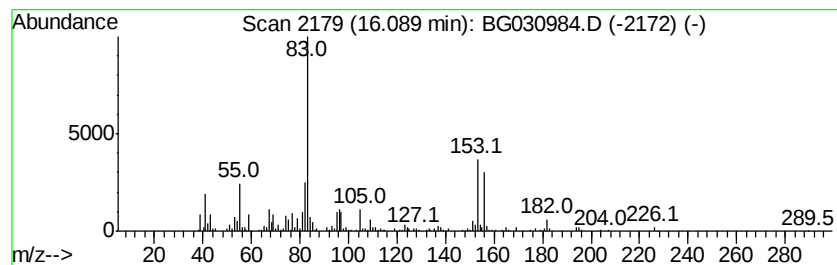
Instrument :
BNA_G
ClientSampleId :
153RD-63RD-ST-DRUM

Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG111017.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 Cyclopentaneacetic acid, 3-... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.		
16.09	11.24 ng	373007	Acenaphthene-d10	14.76		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclopentaneacetic acid, 3-oxo-2...	226	C13H22O3	024851-98-7	93
2		Oxalic acid, cyclohexyl tetradec...	368	C22H40O4	1000309-31-5	38
3		trans-4,4-Dimethyl-2-hexene	112	C8H16	019550-83-5	38
4		Prop-2-ynyl (E)-2-methylbut-2-en...	138	C8H10O2	1000373-72-5	38
5		Cyclohexane, 1,1'-(1,5-pentaned...	236	C17H32	054833-31-7	38



Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA G\DATA\BG111317\
Data File : BG030984.D
Acq On : 13 Nov 2017 22:02
Operator : SJ/JU
Sample : I6436-03
Misc :
ALS Vial : 17 Sample Multiplier: 1

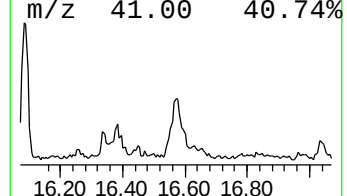
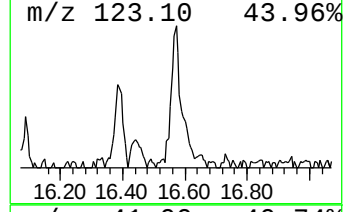
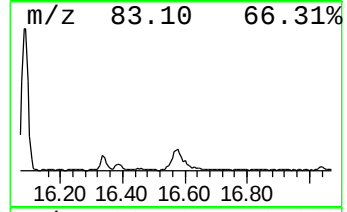
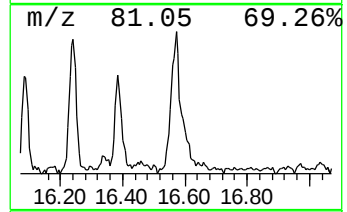
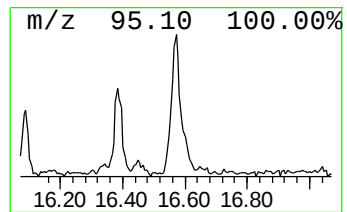
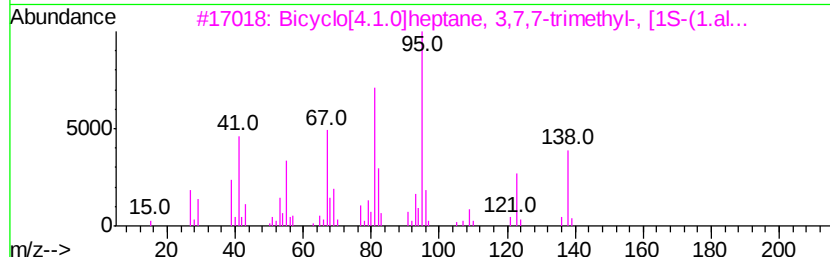
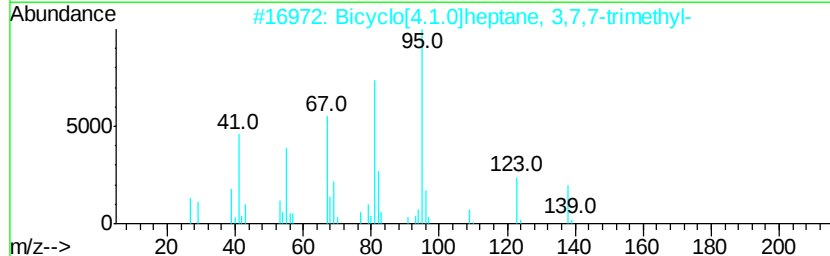
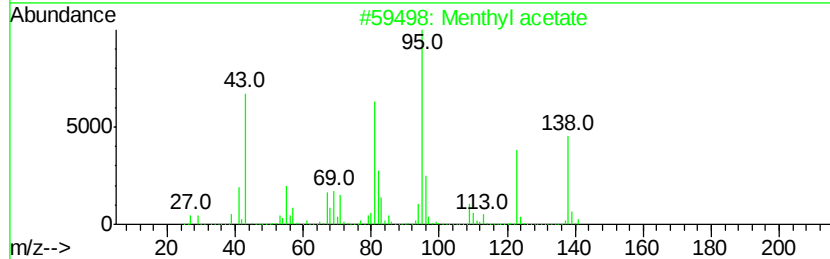
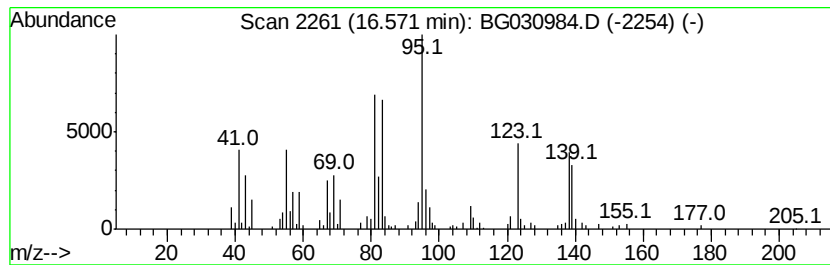
Instrument :
BNA_G
ClientSampleId :
153RD-63RD-ST-DRUM

Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG111017.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 19 Menthyl acetate Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.57	4.25 ng	191263	Phenanthrene-d10	17.49
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Menthyl acetate	198 C12H22O2	000089-48-5 64
2		Bicyclo[4.1.0]heptane, 3,7,7-tri...	138 C10H18	000554-59-6 58
3		Bicyclo[4.1.0]heptane, 3,7,7-tri...	138 C10H18	002778-68-9 53
4		Cyclohexene, 4-methyl-1-(1-methy...	138 C10H18	000500-00-5 53
5		Cyclohexane, 1-methyl-4-(1-methy...	138 C10H18	001124-27-2 53



Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA G\DATA\BG111317\
Data File : BG030984.D
Acq On : 13 Nov 2017 22:02
Operator : SJ/JU
Sample : I6436-03
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
153RD-63RD-ST-DRUM

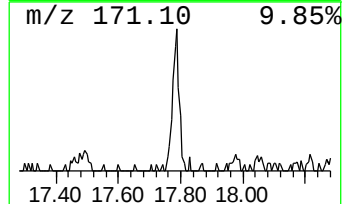
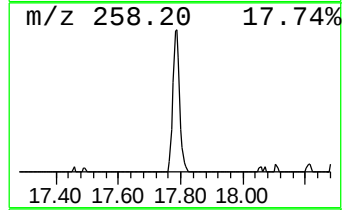
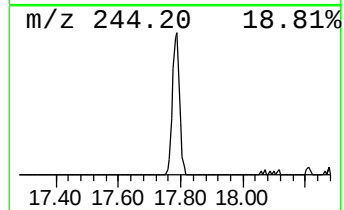
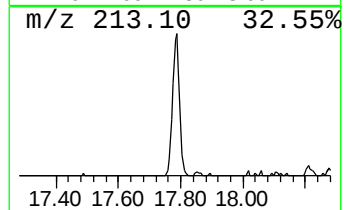
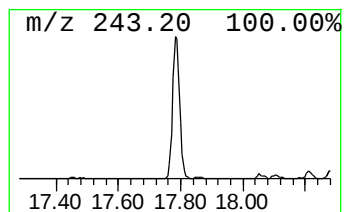
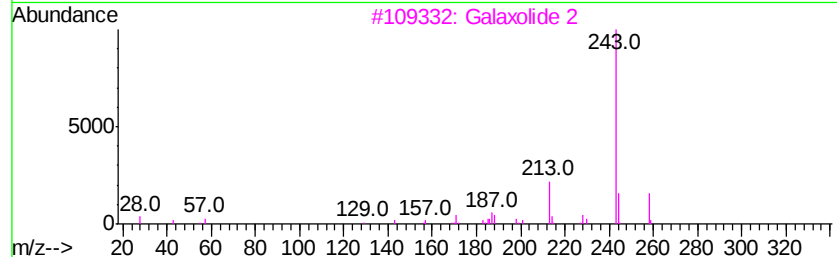
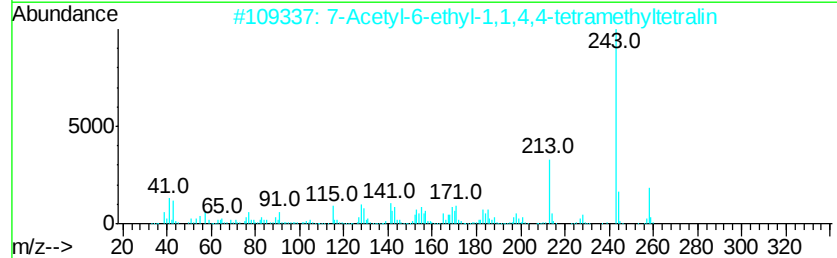
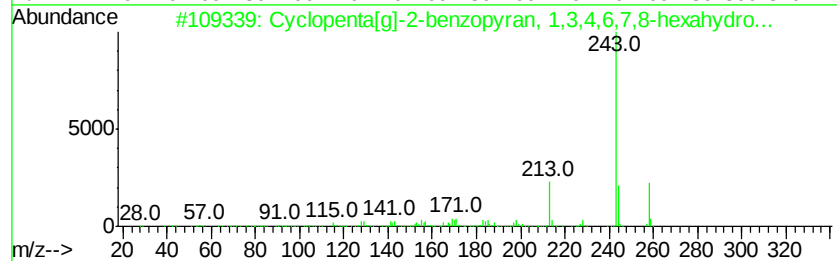
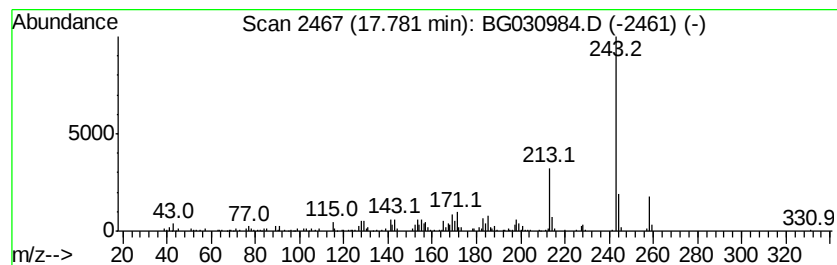
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 20 Cyclopenta[g]-2-benzopyran,... Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.78	4.63 ng	208395	Phenanthrene-d10	17.49

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopenta[g]-2-benzopyran, 1,3,...	258	C18H26O	001222-05-5	97
2			7-Acetyl-6-ethyl-1,1,4,4-tetrame...	258	C18H26O	000088-29-9	96
3			Galaxolide 2	258	C18H26O	1000285-26-7	93
4			Galaxolide 1	258	C18H26O	1000285-26-6	87
5			9-Amino-1-phenyl-3,6-diazahomoad...	243	C15H21N3	147084-69-3	58



Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_G\DATA\BG111317\
 Data File : BG030984.D
 Acq On : 13 Nov 2017 22:02
 Operator : SJ/JU
 Sample : I6436-03
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 153RD-63RD-ST-DRUM

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG111017.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...	5.22	7.7	ng	118074	1	8.14	305153	20.0
1-Methoxy-2-propy...	5.52	10.2	ng	155070	1	8.14	305153	20.0
Benzene, 1,3-dime...	6.14	285.4	ng	4354550	1	8.14	305153	20.0
Benzene, 1-ethyl-...	7.22	18.1	ng	276677	1	8.14	305153	20.0
unknown7.68	7.68	83.2	ng	1268830	1	8.14	305153	20.0
Benzene, 1,2,3-tr...	7.79	31.4	ng	478788	1	8.14	305153	20.0
Benzene, 1-ethyl-...	8.78	5.7	ng	87047	1	8.14	305153	20.0
Hexanoic acid, 3,...	9.78	18.4	ng	384809	2	10.95	419123	20.0
Benzoic acid, 2-m...	11.48	12.7	ng	266551	2	10.95	419123	20.0
Benzoic acid, 4-m...	11.76	7.2	ng	151794	2	10.95	419123	20.0
unknown12.56	12.56	9.8	ng	205462	2	10.95	419123	20.0
Cyclopentaneaceti...	16.09	11.2	ng	373007	3	14.76	663636	20.0
Menthyl acetate	16.57	4.3	ng	191263	4	17.49	900019	20.0
Cyclopenta[g]-2-b...	17.78	4.6	ng	208395	4	17.49	900019	20.0