

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG102716\
 Data File : BG024531.D
 Acq On : 27 Oct 2016 21:30
 Operator : UM/SJ
 Sample : H5432-01
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 GW-1

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-BG102516.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	5.341	420	426	434	rBV	217009	364129	3.27%	0.665%
2	5.805	498	505	514	rBV	712989	1215511	10.92%	2.219%
3	6.310	585	591	600	rBV3	60252	115610	1.04%	0.211%
4	7.403	771	777	788	rBV2	448904	869784	7.81%	1.588%
5	7.797	835	844	851	rBV	1367117	2407740	21.63%	4.396%
6	8.273	917	925	928	rBV2	503807	895168	8.04%	1.634%
7	8.308	928	931	940	rVB2	310323	525218	4.72%	0.959%
8	8.602	971	981	994	rVB	2196997	4059799	36.47%	7.412%
9	8.954	1033	1041	1049	rBV6	60491	138364	1.24%	0.253%
10	9.095	1059	1065	1074	rVB	186462	339159	3.05%	0.619%
11	9.295	1093	1099	1105	rBV4	98121	190950	1.72%	0.349%
12	9.389	1105	1115	1119	rVV4	158805	418549	3.76%	0.764%
13	9.448	1119	1125	1132	rVV	1966559	3647317	32.76%	6.659%
14	9.536	1132	1140	1147	rVB3	204887	450363	4.05%	0.822%
15	9.918	1198	1205	1209	rBV3	86502	218416	1.96%	0.399%
16	10.323	1269	1274	1281	rBV2	66725	113476	1.02%	0.207%
17	10.576	1306	1317	1322	rBV4	220890	541111	4.86%	0.988%
18	10.652	1323	1330	1336	rVB4	181289	348946	3.13%	0.637%
19	10.946	1374	1380	1381	rBV4	82513	129112	1.16%	0.236%
20	11.099	1400	1406	1420	rVB	641475	1379996	12.40%	2.520%
21	11.222	1420	1427	1428	rBV3	219969	338171	3.04%	0.617%
22	11.722	1504	1512	1520	rBV3	162081	384252	3.45%	0.702%
23	12.027	1557	1564	1568	rBV8	48161	113128	1.02%	0.207%
24	12.274	1597	1606	1613	rBV	521135	988882	8.88%	1.806%
25	12.397	1621	1627	1632	rBV6	74020	130680	1.17%	0.239%
26	13.514	1809	1817	1828	rVB	4709186	7460340	67.02%	13.621%
27	13.713	1847	1851	1856	rBV2	81285	133332	1.20%	0.243%
28	14.324	1950	1955	1961	rBV	128151	181840	1.63%	0.332%
29	14.888	2043	2051	2059	rVB	976016	1657342	14.89%	3.026%
30	16.369	2296	2303	2311	rBV	2265528	3508651	31.52%	6.406%
31	16.545	2329	2333	2340	rBV3	67699	121593	1.09%	0.222%
32	17.626	2511	2517	2523	rBV	1246760	2002126	17.99%	3.656%
33	18.472	2657	2661	2668	rBV7	85174	141733	1.27%	0.259%
34	20.018	2919	2924	2933	rBV	1729922	2543387	22.85%	4.644%

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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

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35	20.212	2951	2957	2963	rBV	8519405	11132198	100.00%	20.325%
36	21.921	3242	3248	3258	rVB	1483276	2665844	23.95%	4.867%
37	22.250	3301	3304	3311	rVB10	70272	131854	1.18%	0.241%
38	25.364	3822	3834	3845	rVB2	911875	2765701	24.84%	5.050%

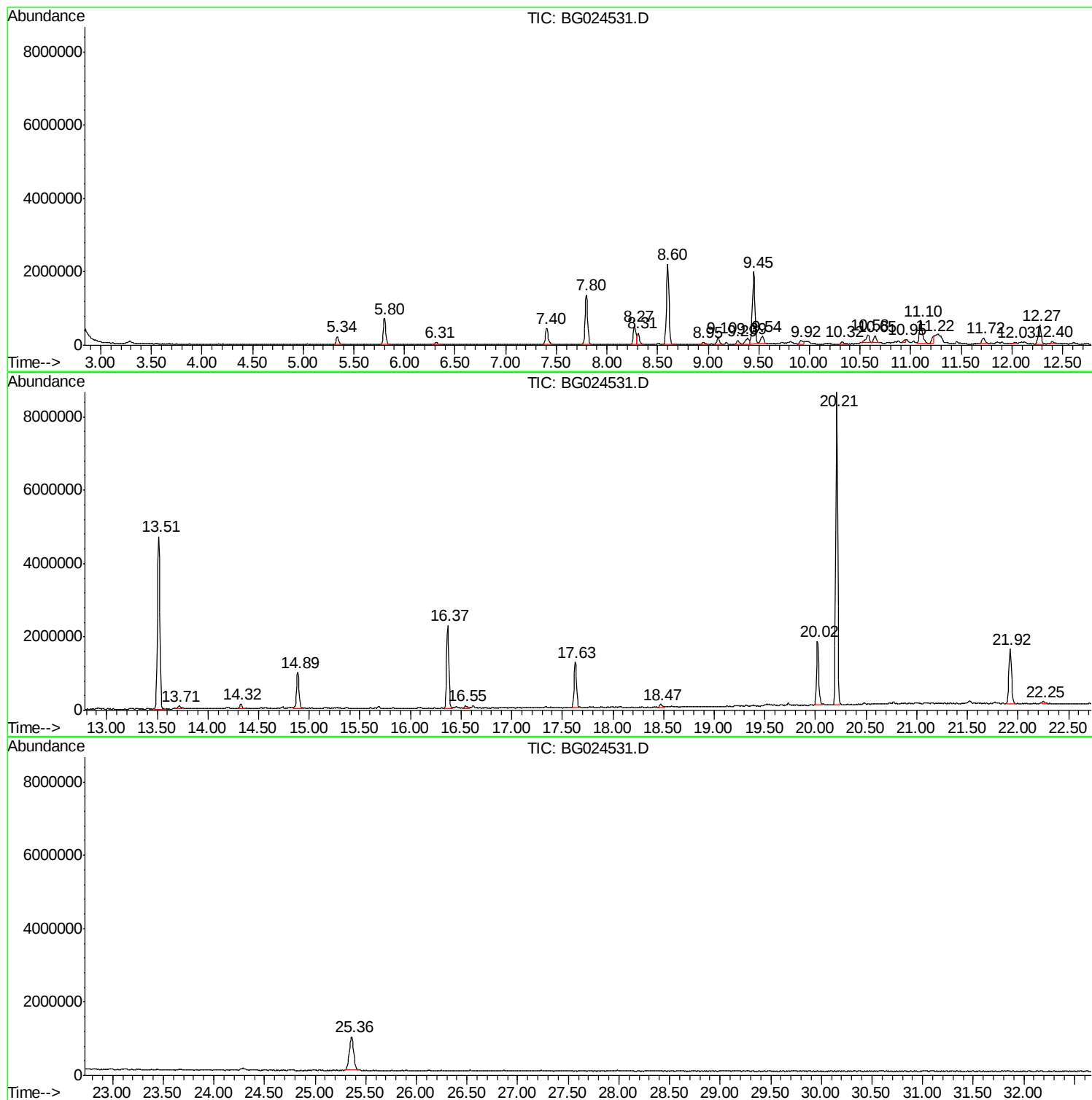
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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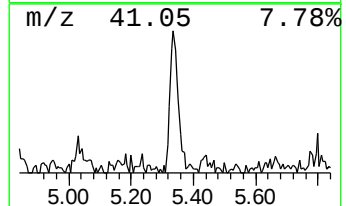
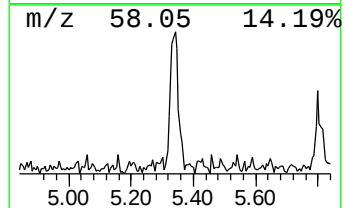
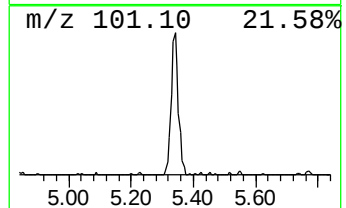
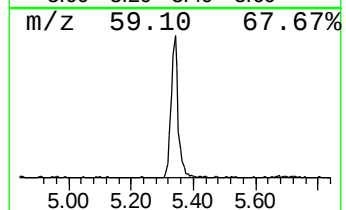
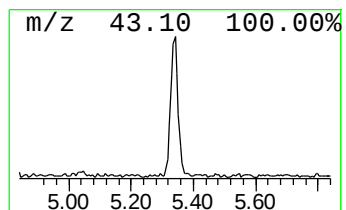
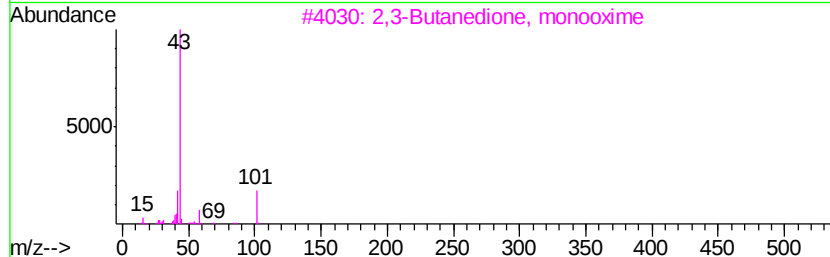
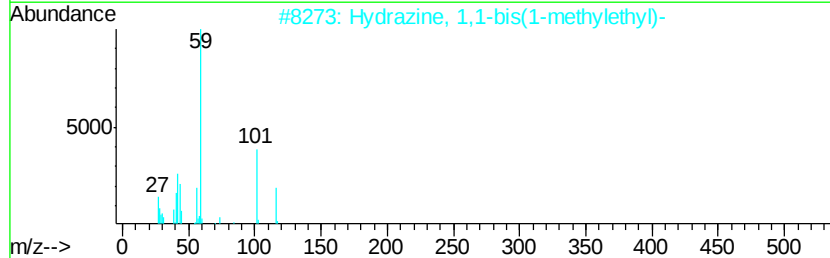
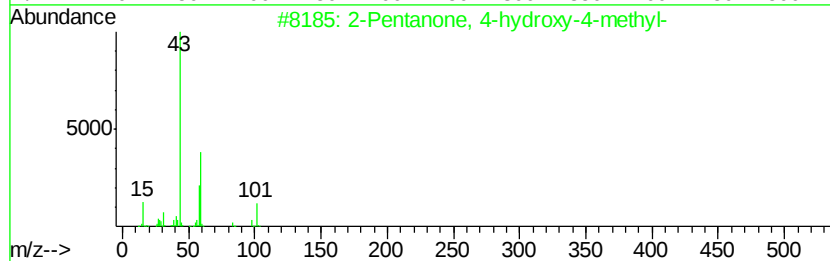
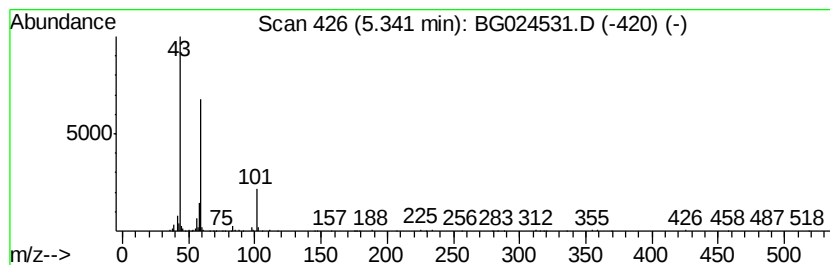
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG102516.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.34	8.14 ng	364129	1,4-Dichlorobenzene-d4	8.27

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2			Hydrazine, 1,1-bis(1-methylethyl)-	116	C6H16N2	000921-14-2	38
3			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	35
4			3-Hexanol	102	C6H14O	000623-37-0	23
5			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	17



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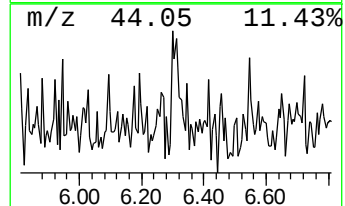
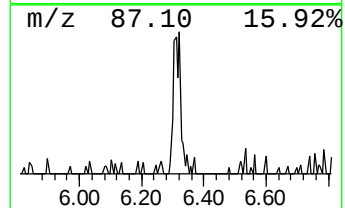
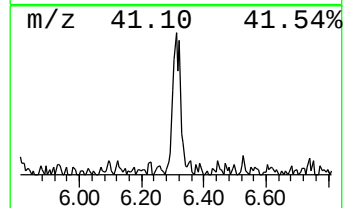
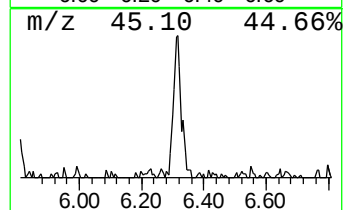
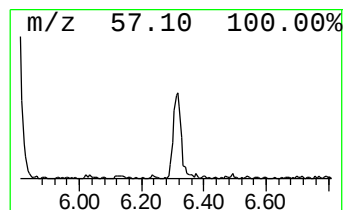
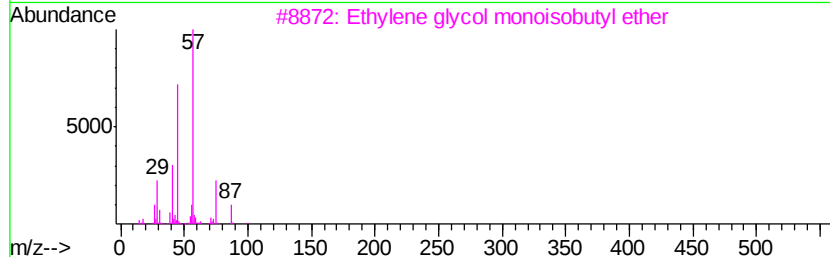
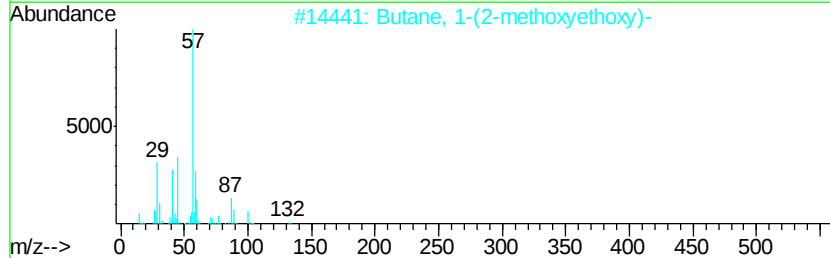
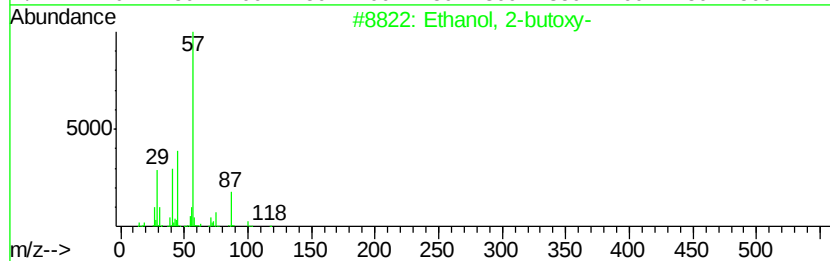
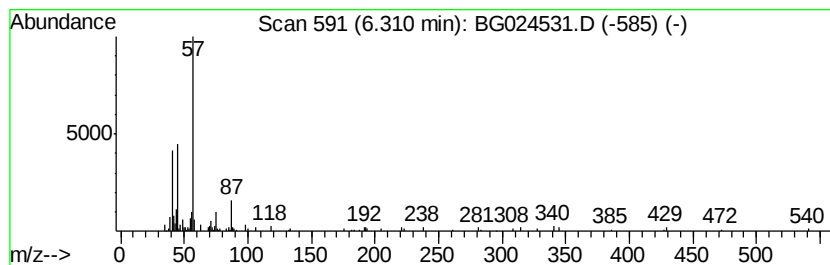
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Ethanol, 2-butoxy- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.31	2.58 ng	115610	1,4-Dichlorobenzene-d4	8.27

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-butoxy-	118	C6H14O2	000111-76-2	59
2			Butane, 1-(2-methoxyethoxy)-	132	C7H16O2	013343-98-1	45
3			Ethylene glycol monoisobutyl ether	118	C6H14O2	004439-24-1	42
4			2-Pentanol, 4,4-dimethyl-	116	C7H16O	006144-93-0	40
5			Butanoic acid, 2-oxo-	102	C4H6O3	000600-18-0	40



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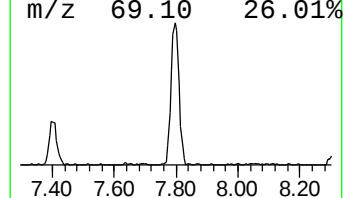
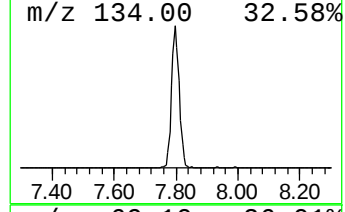
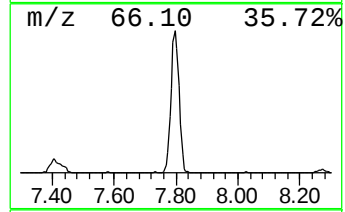
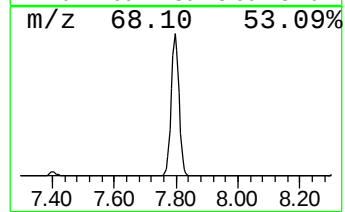
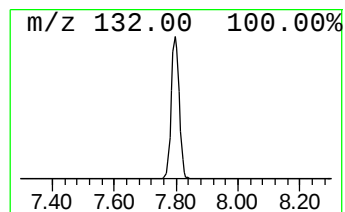
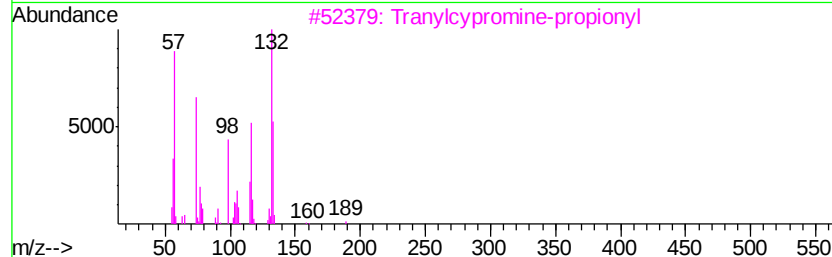
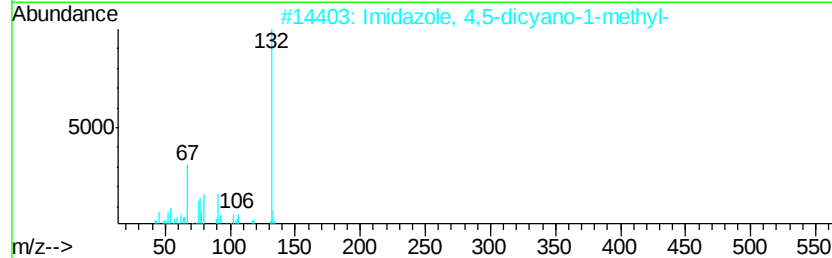
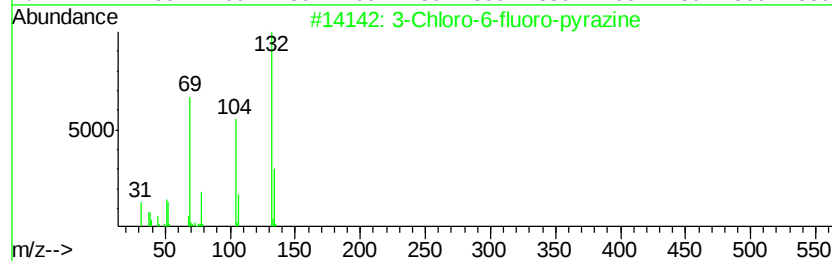
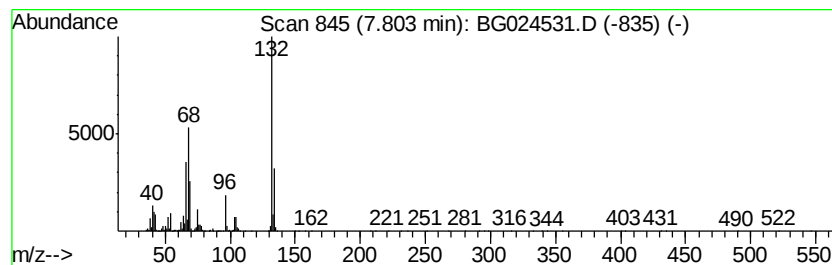
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 unknown7.80 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.80	53.79 ng	2407740	1,4-Dichlorobenzene-d4	8.27

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		Imidazole, 4,5-dicyano-1-methyl-	132	C6H4N4	019485-35-9	25
3		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	22
4		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
5		Cyclopropanamine, 1-phenyl-	133	C9H11N	1000351-26-2	22



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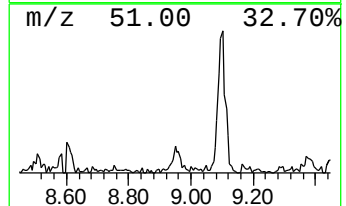
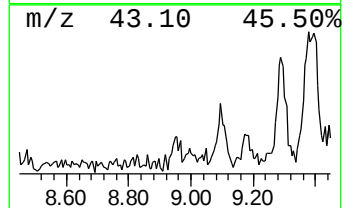
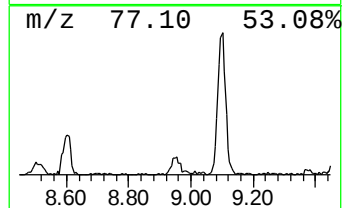
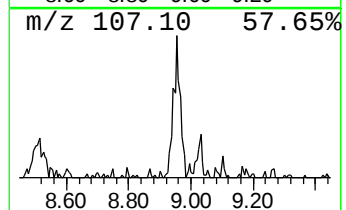
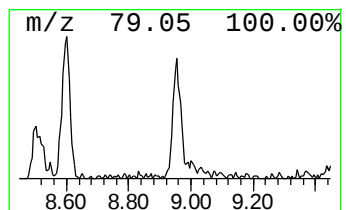
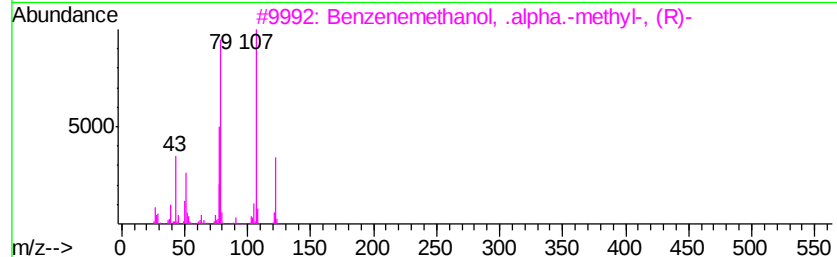
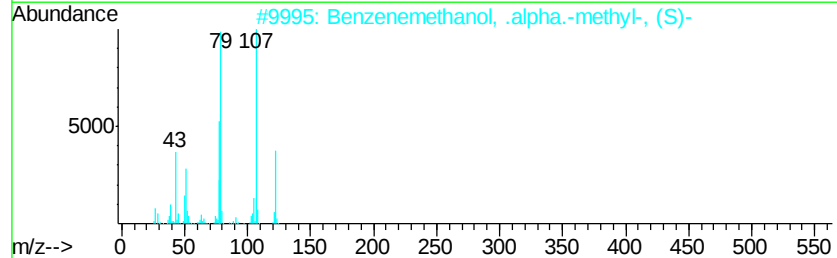
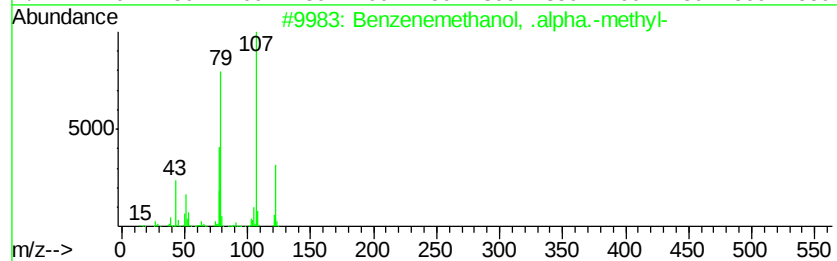
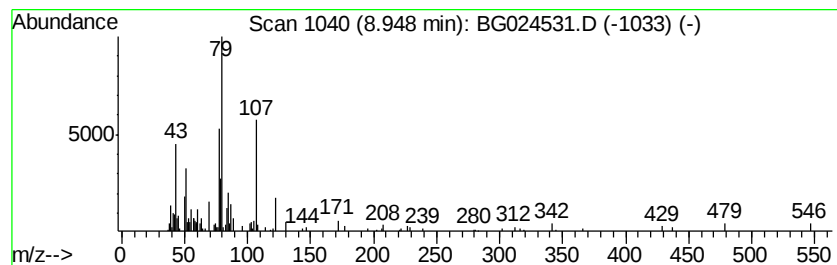
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Benzenemethanol, .alpha.-me... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.95	3.09 ng	138364	1,4-Dichlorobenzene-d4	8.27

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzenemethanol, .alpha.-methyl-	122	C8H10O	000098-85-1	90
2		Benzenemethanol, .alpha.-methyl-...	122	C8H10O	001445-91-6	86
3		Benzenemethanol, .alpha.-methyl-...	122	C8H10O	001517-69-7	86
4		(R)-(+)-1-Phenyl-1-propanol	136	C9H12O	001565-74-8	72
5		1-Hydroxy,1-phenyl-2-propanone	150	C9H10O2	1000222-86-8	72



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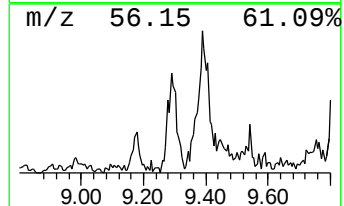
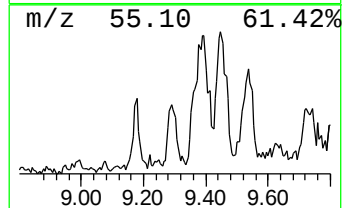
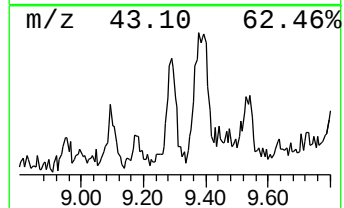
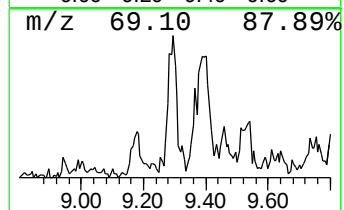
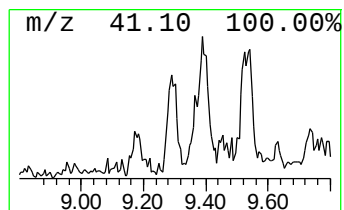
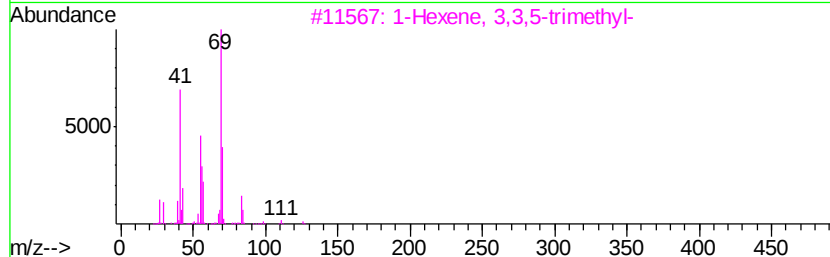
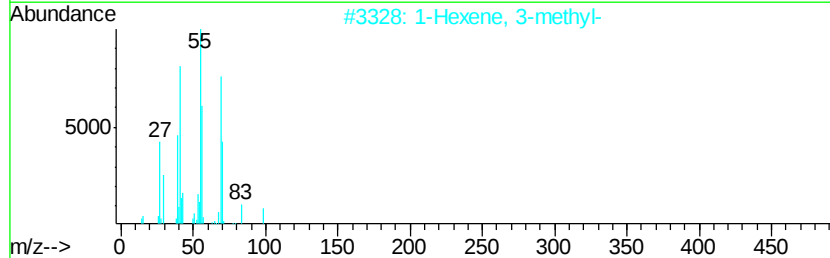
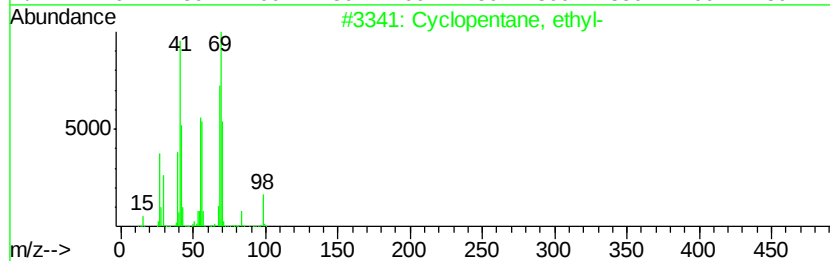
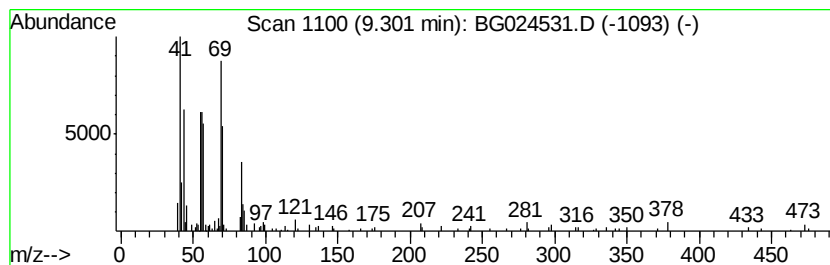
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Cyclopentane, ethyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.30	4.27 ng	190950	1,4-Dichlorobenzene-d4	8.27

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopentane, ethyl-	98	C7H14	001640-89-7	74
2		1-Hexene, 3-methyl-	98	C7H14	003404-61-3	72
3		1-Hexene, 3,3,5-trimethyl-	126	C9H18	013427-43-5	58
4		3-Methyl-2-hexene	98	C7H14	017618-77-8	53
5		1-Pentene, 2,3-dimethyl-	98	C7H14	003404-72-6	50



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG102716\
Data File : BG024531.D
Acq On : 27 Oct 2016 21:30
Operator : UM/SJ
Sample : H5432-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

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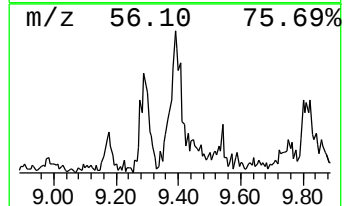
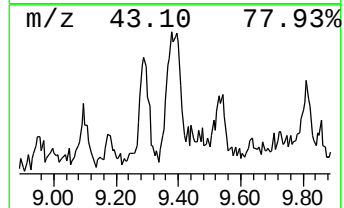
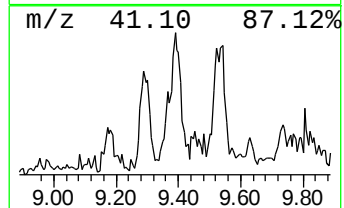
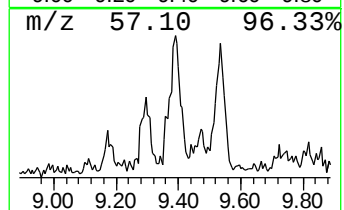
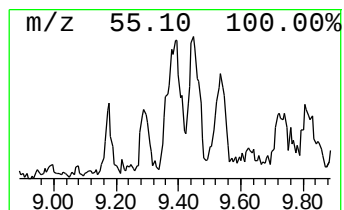
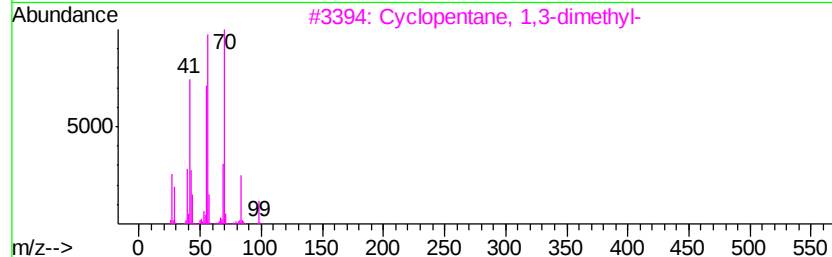
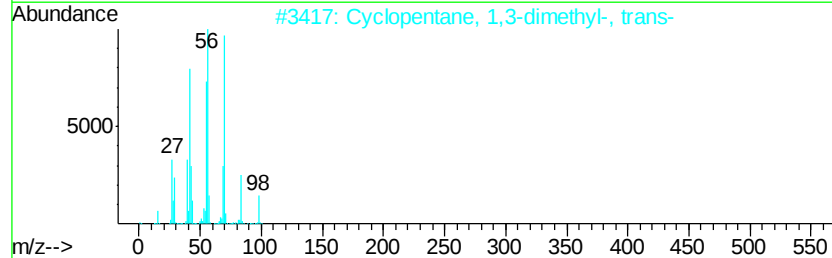
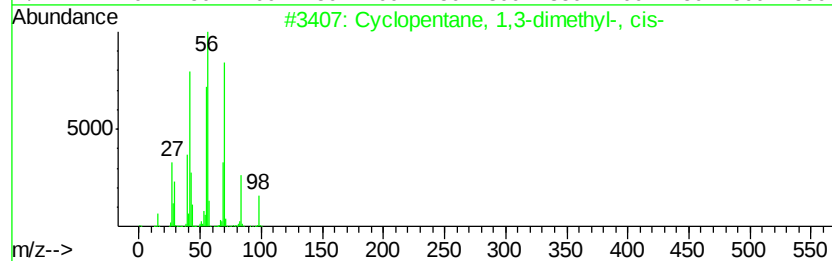
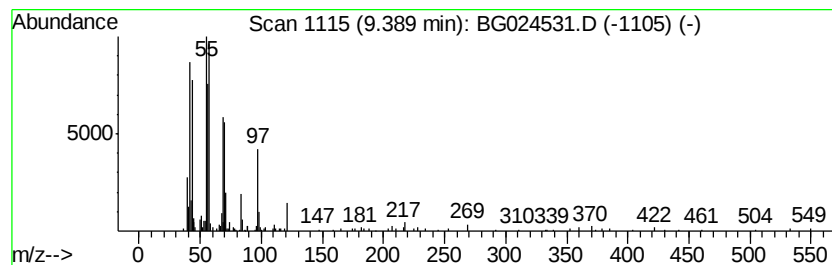
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Cyclopentane, 1,3-dimethyl-... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.39	9.35 ng	418549	1,4-Dichlorobenzene-d4	8.27

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopentane, 1,3-dimethyl-, cis-	98	C7H14	002532-58-3	52
2		Cyclopentane, 1,3-dimethyl-, trans-	98	C7H14	001759-58-6	52
3		Cyclopentane, 1,3-dimethyl-	98	C7H14	002453-00-1	52
4		1-Hexene, 3-methyl-	98	C7H14	003404-61-3	50
5		Cyclopropane, 1-methyl-2-(3-meth...	140	C10H20	062238-07-7	47



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG102716\
Data File : BG024531.D
Acq On : 27 Oct 2016 21:30
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Misc :
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ClientSampleId :
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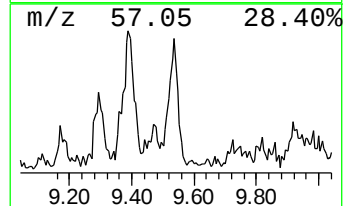
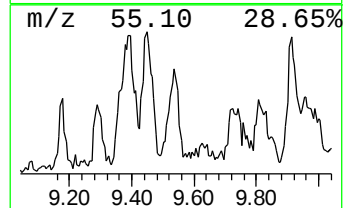
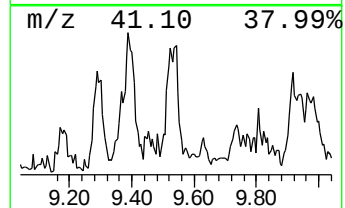
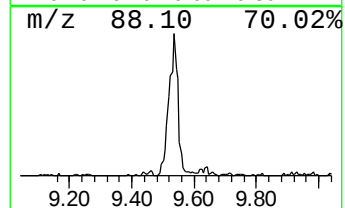
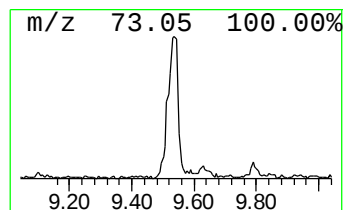
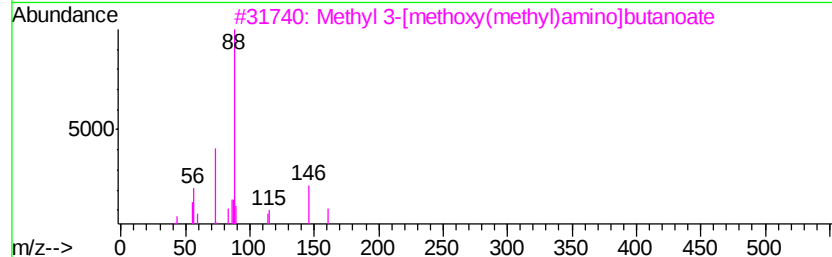
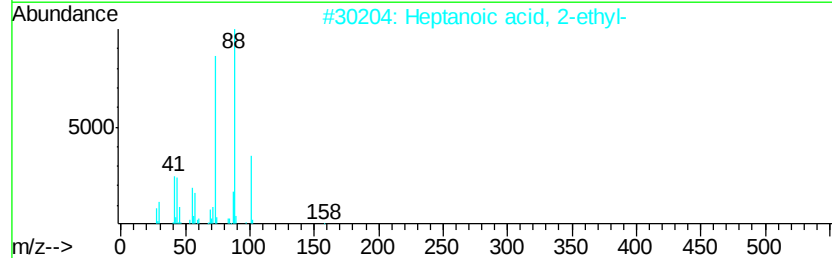
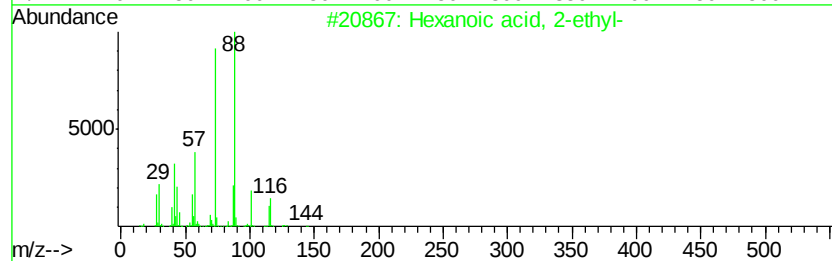
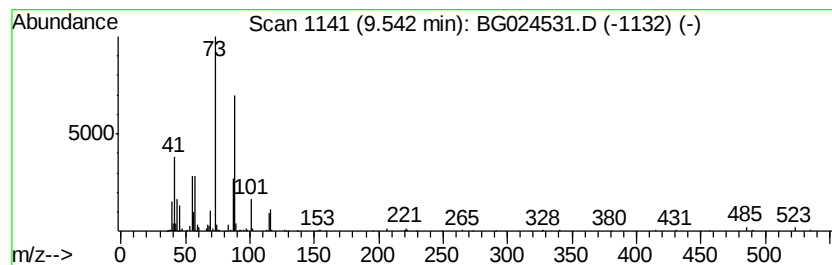
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Hexanoic acid, 2-ethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.54	10.06 ng	450363	1,4-Dichlorobenzene-d4	8.27

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	56
2		Heptanoic acid, 2-ethyl-	158	C9H18O2	003274-29-1	53
3		Methyl 3-[methoxy(methyl)amino]b...	161	C7H15N03	959080-02-5	36
4		Acetic acid, diethyl-	116	C6H12O2	000088-09-5	33
5		.alpha.-Ethyl aspartate	161	C6H11N04	1000132-78-1	23



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG102716\
Data File : BG024531.D
Acq On : 27 Oct 2016 21:30
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Sample : H5432-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

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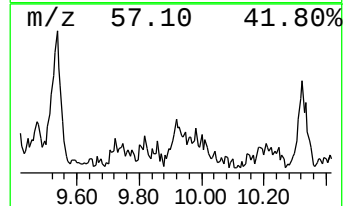
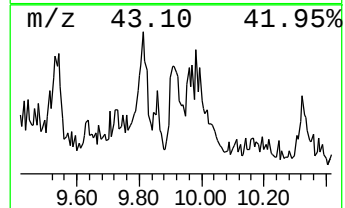
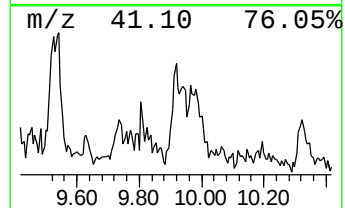
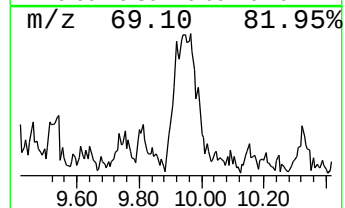
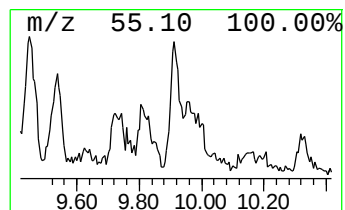
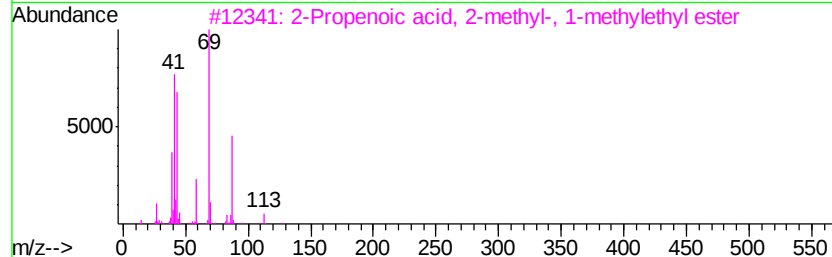
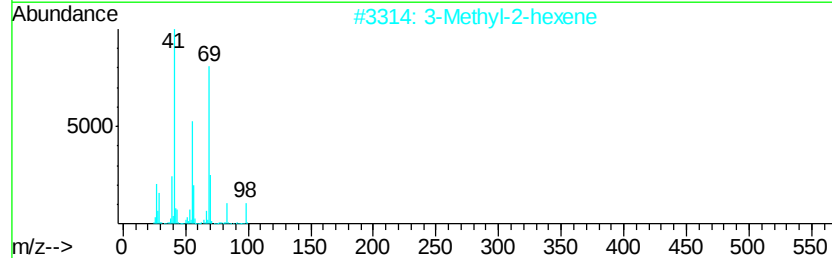
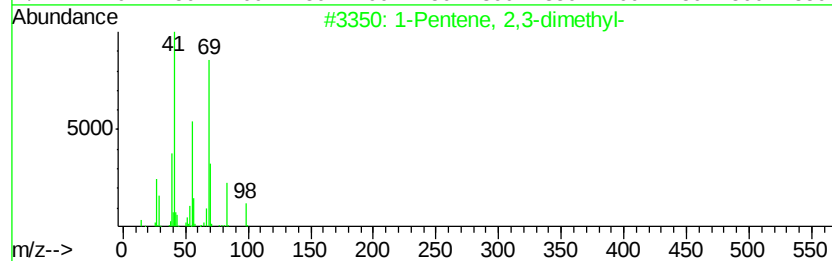
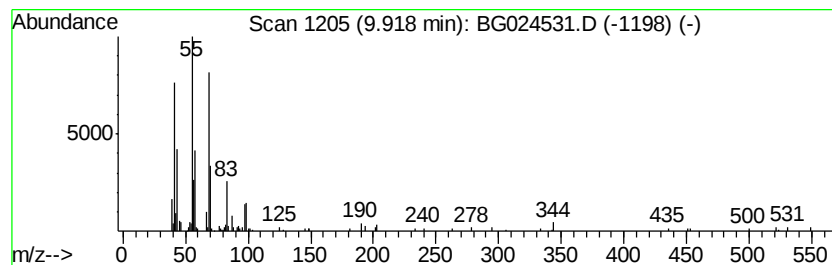
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 1-Pentene, 2,3-dimethyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.92	3.17 ng	218416	Napthalene-d8	11.10

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Pentene, 2,3-dimethyl-	98	C7H14	003404-72-6	64
2		3-Methyl-2-hexene	98	C7H14	017618-77-8	59
3		2-Propenoic acid, 2-methyl-, 1-m...	128	C7H12O2	004655-34-9	45
4		(Z)-4-Methyl-2-hexene	98	C7H14	003683-19-0	45
5		6-Dodecene, (E)-	168	C12H24	007206-17-9	45



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG102716\
Data File : BG024531.D
Acq On : 27 Oct 2016 21:30
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Misc :
ALS Vial : 8 Sample Multiplier: 1

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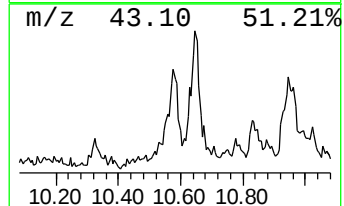
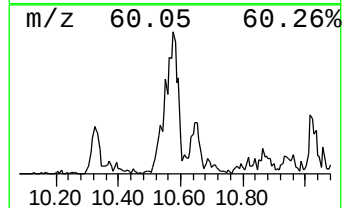
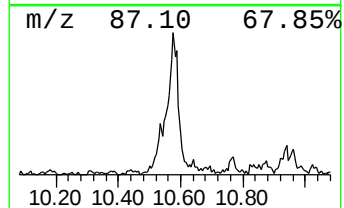
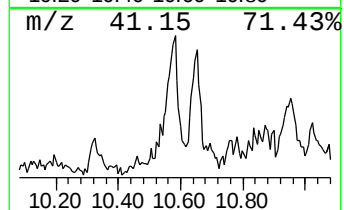
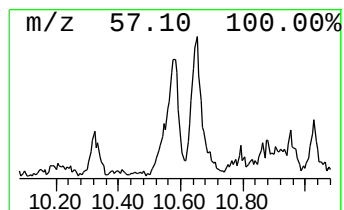
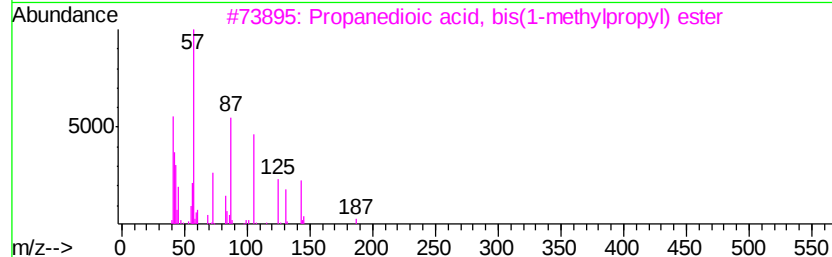
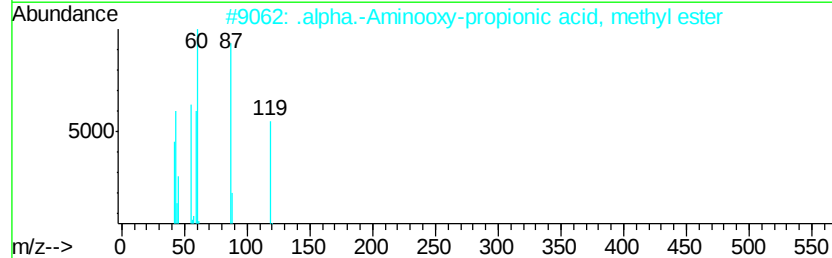
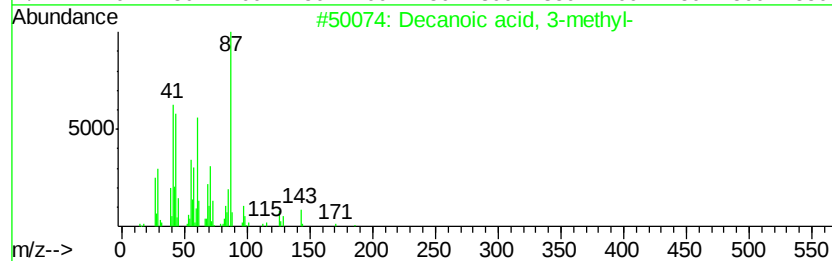
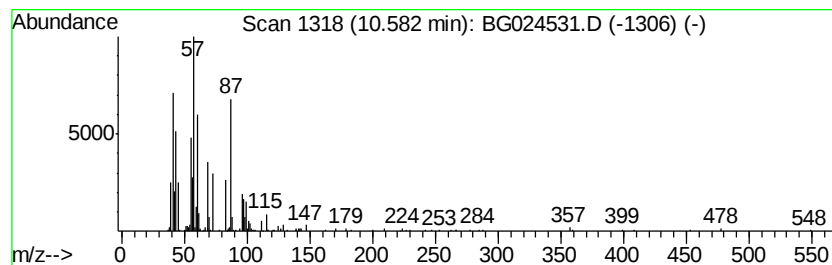
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Decanoic acid, 3-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.58	7.84 ng	541111	Naphthalene-d8	11.10

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Decanoic acid, 3-methyl-	186	C11H22O2	060308-82-9	50
2		.alpha.-Aminooxy-propionic acid,...	119	C4H9NO3	091666-48-7	32
3		Propanedioic acid, bis(1-methylp...	216	C11H20O4	032260-07-4	32
4		Thiophene, tetrahydro-2-methyl-	102	C5H10S	001795-09-1	27
5		2,7-Dimethyl-4,6-dioxanonane	160	C9H20O2	1000334-82-5	27



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG102716\
Data File : BG024531.D
Acq On : 27 Oct 2016 21:30
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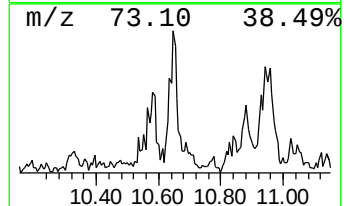
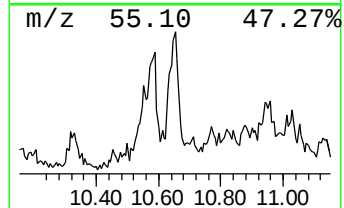
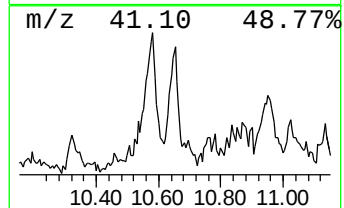
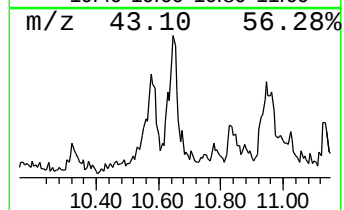
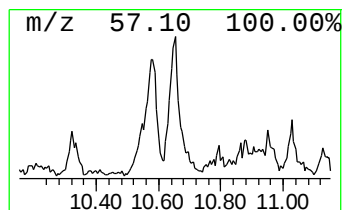
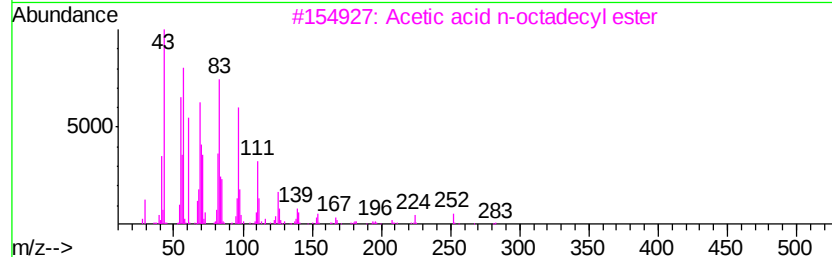
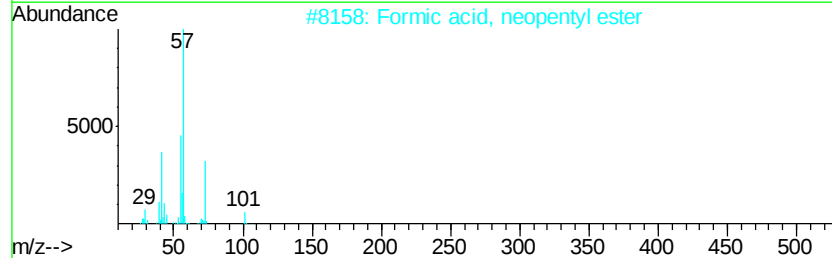
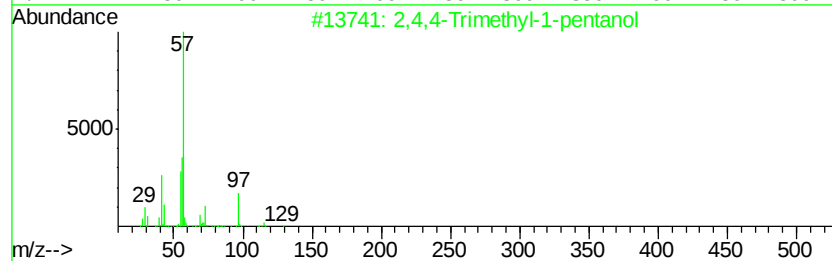
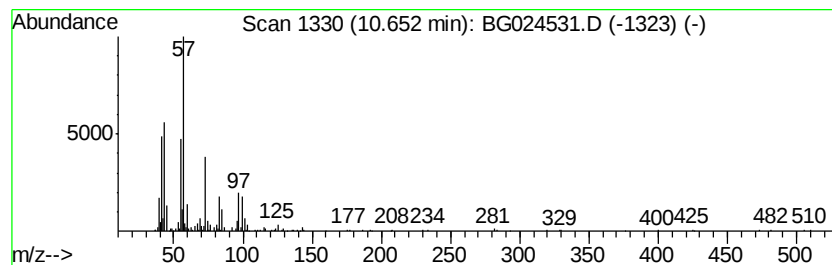
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 unknown10.65 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.65	5.06 ng	348946	Naphthalene-d8	11.10

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,4,4-Trimethyl-1-pentanol			130	C8H18O	016325-63-6	43
2	Formic acid, neopentyl ester			116	C6H12O2	1000367-90-3	35
3	Acetic acid n-octadecyl ester			312	C20H40O2	000822-23-1	22
4	Hexane, 4-ethyl-2-methyl-			128	C9H20	003074-75-7	22
5	Sulfurous acid, butyl 2-propyl e...			180	C7H16O3S	1000309-11-4	22



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG102716\
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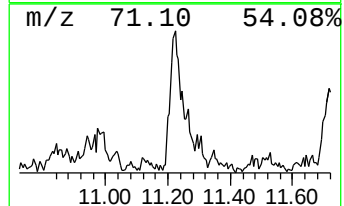
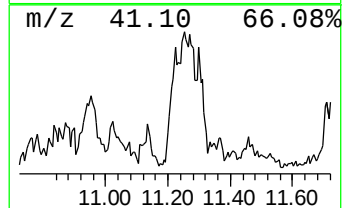
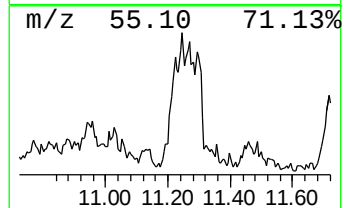
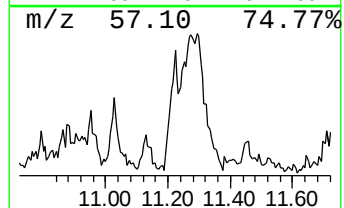
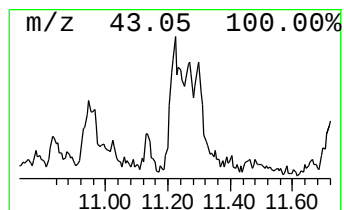
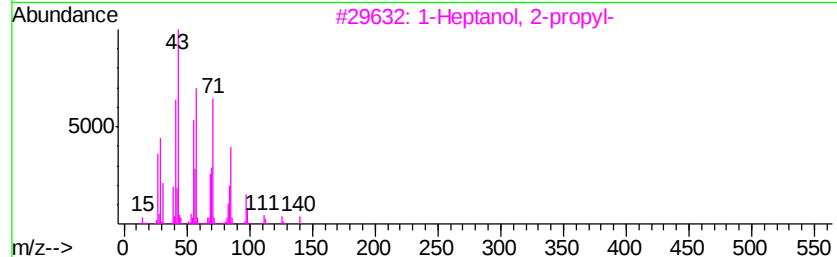
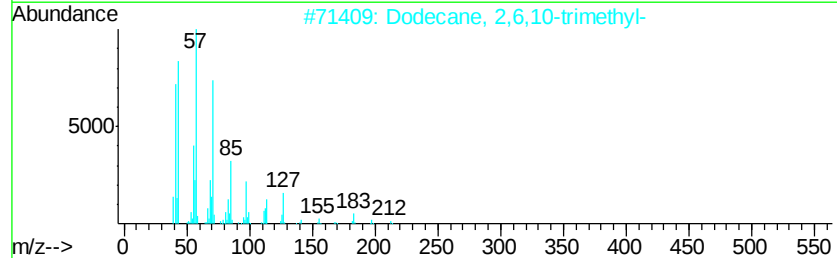
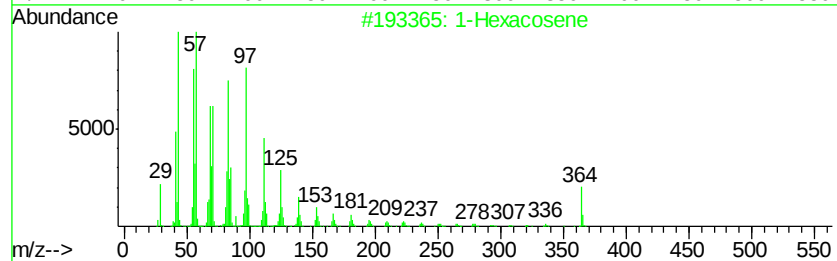
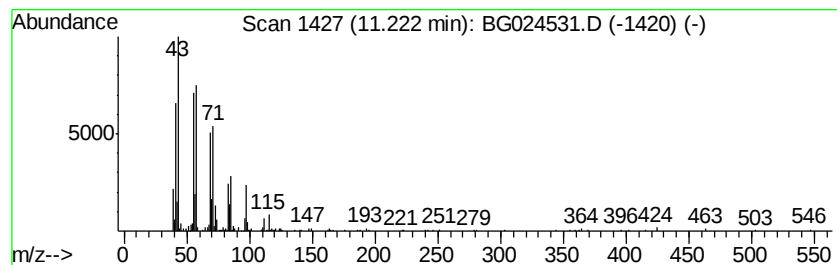
Instrument :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 1-Hexacosene Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.22	4.90 ng	338171	Naphthalene-d8	11.10
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	1-Hexacosene		364 C26H52	018835-33-1 60
2	Dodecane, 2,6,10-trimethyl-		212 C15H32	003891-98-3 50
3	1-Heptanol, 2-propyl-		158 C10H22O	010042-59-8 47
4	1-Decanol, 2-hexyl-		242 C16H34O	002425-77-6 47
5	Tetradecane, 2,6,10-trimethyl-		240 C17H36	014905-56-7 46



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG102716\
Data File : BG024531.D
Acq On : 27 Oct 2016 21:30
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Sample : H5432-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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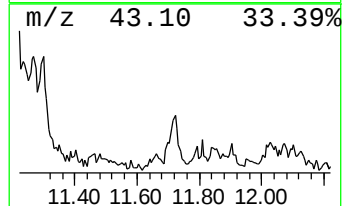
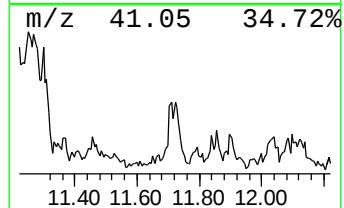
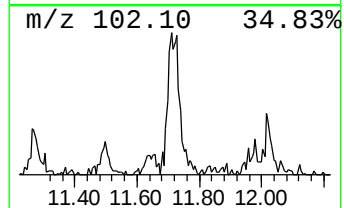
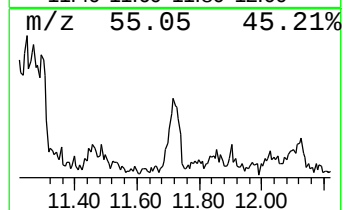
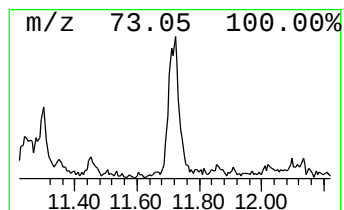
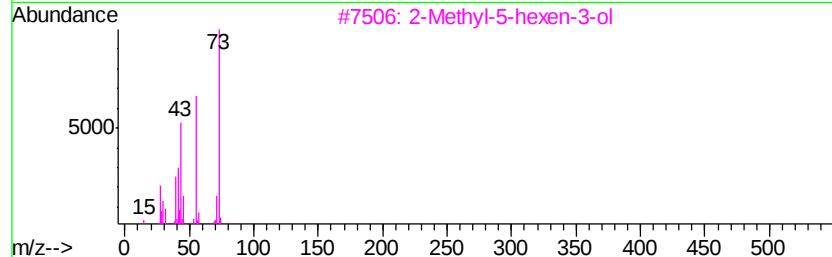
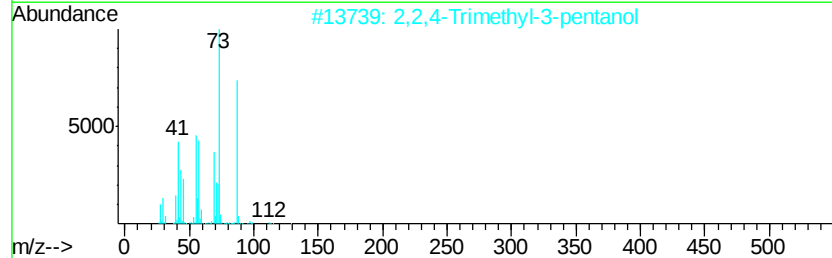
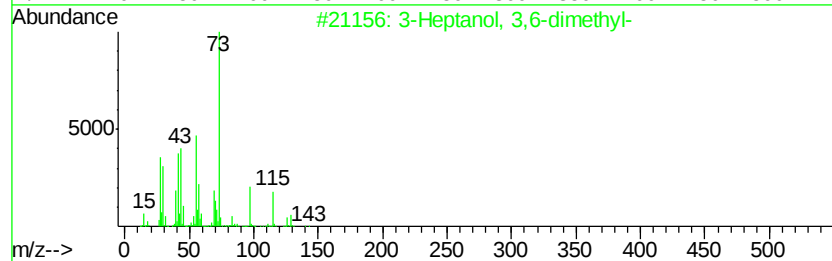
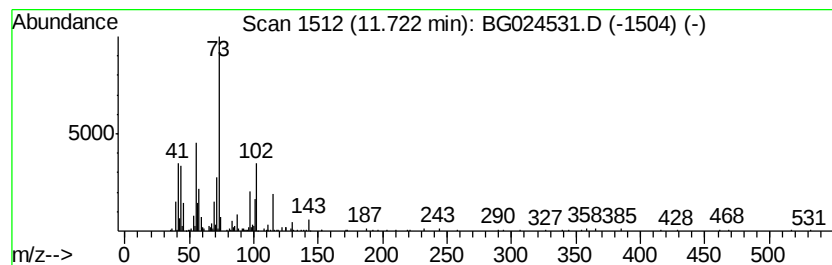
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG102516.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 unknown11.72 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.72	5.57 ng	384252	Naphthalene-d8	11.10

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Heptanol, 3,6-dimethyl-	144	C9H20O	001573-28-0	27
2			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	16
3			2-Methyl-5-hexen-3-ol	114	C7H14O	032815-70-6	14
4			3,4-Dimethyl-5-hexen-3-ol	128	C8H16O	001569-45-5	10
5			3,4,4-Trimethyl-3-pentanol	130	C8H18O	007294-05-5	10



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG102716\
Data File : BG024531.D
Acq On : 27 Oct 2016 21:30
Operator : UM/SJ
Sample : H5432-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

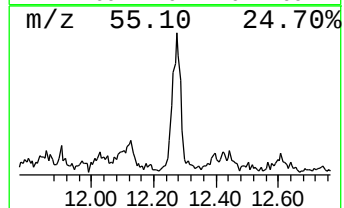
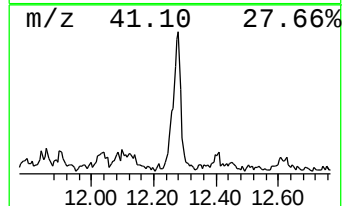
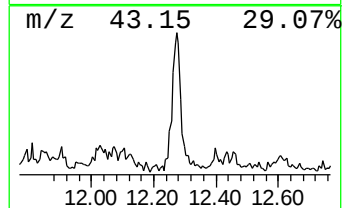
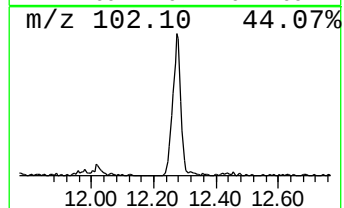
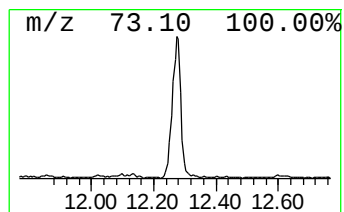
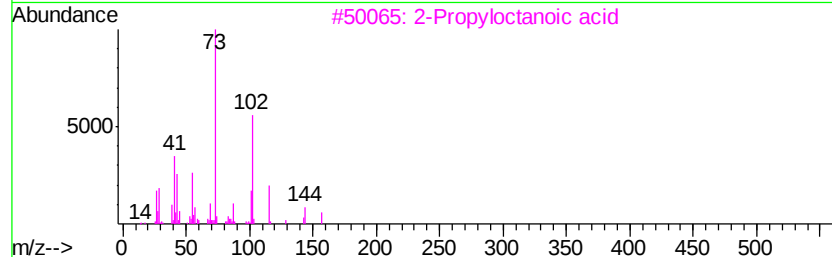
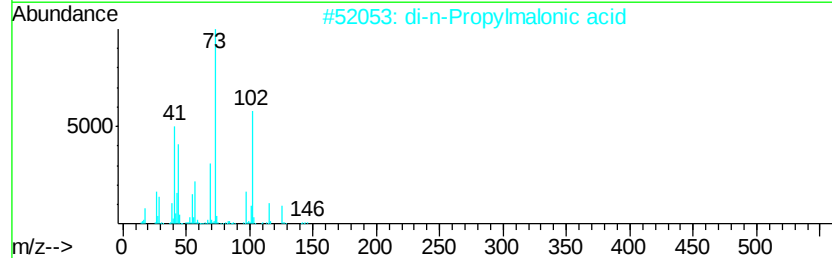
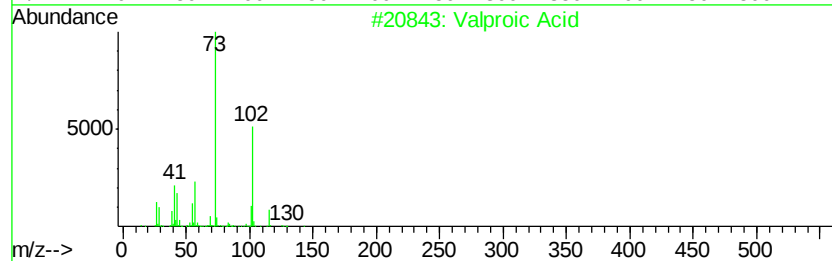
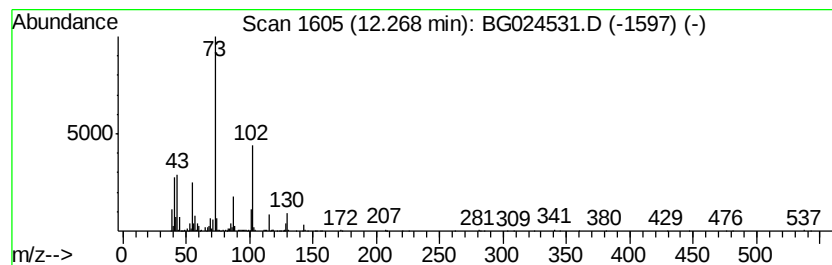
Instrument :
BNA_G
ClientSampleId :
GW-1

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG102516.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Valproic Acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.27	14.33 ng	988882	Napthalene-d8	11.10
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Valproic Acid	144 C8H16O2	000099-66-1	53
2	di-n-Propylmalonic acid	188 C9H16O4	001636-27-7	42
3	2-Propyloctanoic acid	186 C11H22O2	031080-41-8	40
4	Butane, 2-methoxy-2-methyl-	102 C6H14O	000994-05-8	10
5	l-Inositol	180 C6H12O6	1000007-49-1	9



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG102716\
Data File : BG024531.D
Acq On : 27 Oct 2016 21:30
Operator : UM/SJ
Sample : H5432-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
GW-1

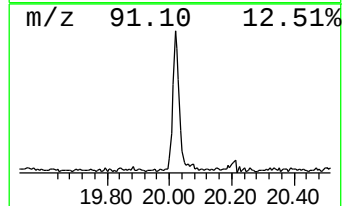
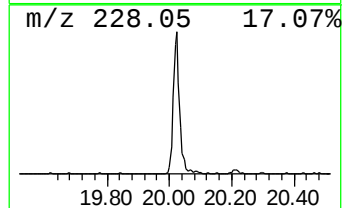
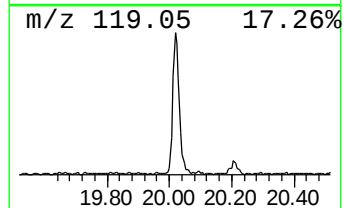
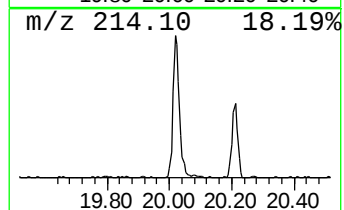
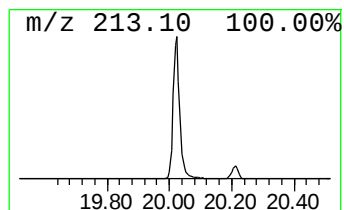
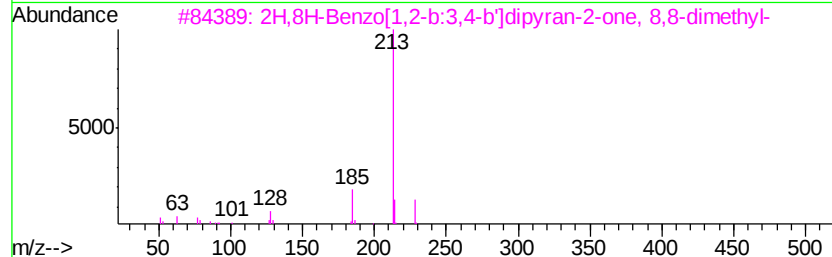
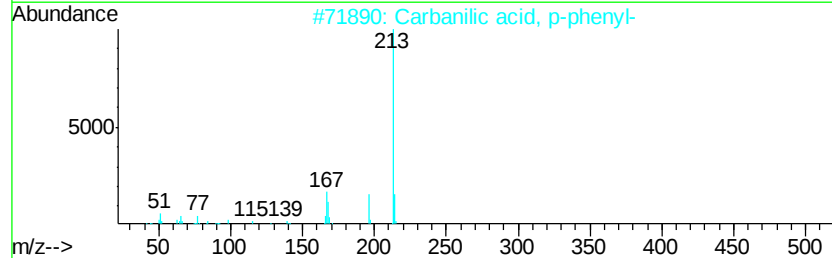
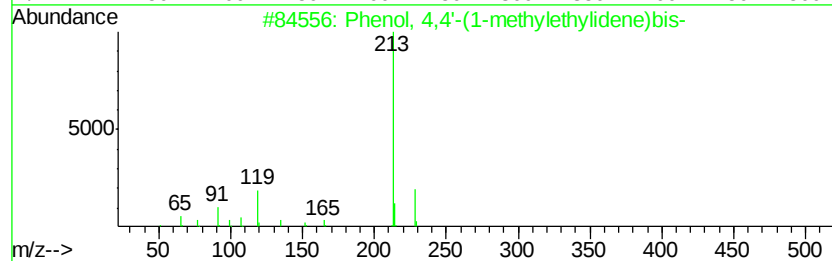
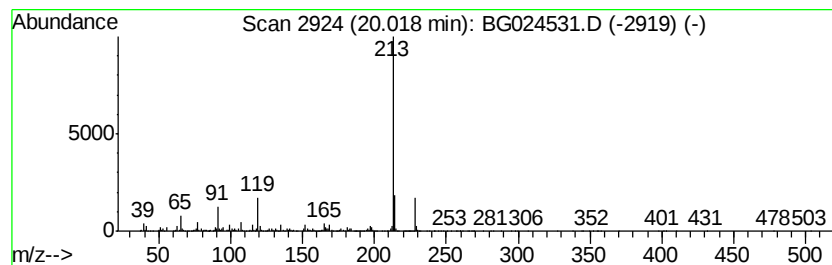
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG102516.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Phenol, 4,4'-(1-methylethyl... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.02	19.08 ng	2543390	Chrysene-d12	21.92

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phenol, 4,4'-(1-methylethylidene...	228	C15H16O2	000080-05-7	96
2		Carbanilic acid, p-phenyl-	213	C13H11NO2	004474-53-7	50
3		2H,8H-Benzo[1,2-b:3,4-b']dipyran...	228	C14H12O3	000523-59-1	45
4		2,5-bis(Trimethylsilyl)thiophene	228	C10H20SSi2	017906-71-7	42
5		Indole,6-methyl-2-(2-thienyl)-	213	C13H11NS	296245-70-0	40



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG102716\
 Data File : BG024531.D
 Acq On : 27 Oct 2016 21:30
 Operator : UM/SJ
 Sample : H5432-01
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 GW-1

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG102516.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...	5.34	8.1 ng		364129	1	8.27	895168	20.0
Ethanol, 2-butoxy-	6.31	2.6 ng		115610	1	8.27	895168	20.0
unknown7.80	7.80	53.8 ng		2407740	1	8.27	895168	20.0
Benzenemethanol, ...	8.95	3.1 ng		138364	1	8.27	895168	20.0
Cyclopentane, ethyl-	9.30	4.3 ng		190950	1	8.27	895168	20.0
Cyclopentane, 1,3...	9.39	9.3 ng		418549	1	8.27	895168	20.0
Hexanoic acid, 2-...	9.54	10.1 ng		450363	1	8.27	895168	20.0
1-Pentene, 2,3-di...	9.92	3.2 ng		218416	2	11.10	1380000	20.0
Decanoic acid, 3-...	10.58	7.8 ng		541111	2	11.10	1380000	20.0
unknown10.65	10.65	5.1 ng		348946	2	11.10	1380000	20.0
1-Hexacosene	11.22	4.9 ng		338171	2	11.10	1380000	20.0
unknown11.72	11.72	5.6 ng		384252	2	11.10	1380000	20.0
Valproic Acid	12.27	14.3 ng		988882	2	11.10	1380000	20.0
Phenol, 4,4'-(1-m...	20.02	19.1 ng		2543390	5	21.92	2665840	20.0