

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF040915\
 Data File : BF078385.D
 Acq On : 10 Apr 2015 5:54
 Operator : TP/IZ
 Sample : G1782-05
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB10(2)

Integration Parameters: rteint.p
 Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0

Filtering: 5
 Min Area: 3 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF040115.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	0.990	7	9	13	rVB	38967	58215	2.86%	0.440%
2	1.058	13	15	18	rVV	649197	633566	31.18%	4.791%
3	1.173	22	25	32	rVB	84171	162260	7.98%	1.227%
4	1.310	35	37	40	rVB	265999	324697	15.98%	2.455%
5	1.538	55	57	64	rVV	90537	159205	7.83%	1.204%
6	1.630	64	65	101	rVB	859423	2032252	100.00%	15.367%
7	4.510	315	317	320	rBV	24179	33302	1.64%	0.252%
8	4.567	320	322	331	rVB	59411	82926	4.08%	0.627%
9	5.150	371	373	376	rBV	632313	729357	35.89%	5.515%
10	6.487	488	490	498	rVB	792377	786282	38.69%	5.946%
11	6.602	498	500	503	rBV	835239	811386	39.93%	6.135%
12	6.887	523	525	527	rBV	182269	196116	9.65%	1.483%
13	7.070	539	541	543	rBV	687032	623022	30.66%	4.711%
14	7.585	584	586	588	rBV	525886	455001	22.39%	3.441%
15	8.465	660	663	665	rBV	281423	278501	13.70%	2.106%
16	9.813	778	781	783	rBV	1067623	1021474	50.26%	7.724%
17	10.328	824	826	828	rBV	29978	34303	1.69%	0.259%
18	10.625	849	852	854	rBV	324088	346709	17.06%	2.622%
19	11.528	926	931	933	rVB	23248	22884	1.13%	0.173%
20	11.597	935	937	940	rBV	554354	547779	26.95%	4.142%
21	11.848	957	959	960	rBV	25096	27179	1.34%	0.206%
22	12.134	982	984	987	rBV	20926	30601	1.51%	0.231%
23	12.442	1009	1011	1013	rBV	227057	316803	15.59%	2.396%
24	13.197	1075	1077	1080	rVV	101980	182809	9.00%	1.382%
25	13.277	1082	1084	1087	rVV2	23312	48338	2.38%	0.366%
26	13.334	1087	1089	1092	rVV	15243	21976	1.08%	0.166%
27	13.414	1092	1096	1098	rVV4	13293	37462	1.84%	0.283%
28	13.802	1127	1130	1131	rBV3	14709	22108	1.09%	0.167%
29	13.917	1137	1140	1141	rBV3	20646	35208	1.73%	0.266%
30	13.940	1141	1142	1144	rVB	23050	20542	1.01%	0.155%
31	14.100	1147	1156	1161	rBV	130347	481201	23.68%	3.639%
32	14.180	1161	1163	1165	rVV3	28326	61772	3.04%	0.467%
33	14.214	1165	1166	1171	rVV4	29780	52571	2.59%	0.398%
34	14.294	1171	1173	1177	rVB2	12655	22063	1.09%	0.167%

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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

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35	14.443	1183	1186	1188	rBV	1122524	1149683	56.57%	8.693%
36	14.763	1212	1214	1218	rBV5	17850	30911	1.52%	0.234%
37	15.174	1249	1250	1253	rBV3	14968	29729	1.46%	0.225%
38	15.631	1287	1290	1294	rBV	165243	268987	13.24%	2.034%
39	15.711	1294	1297	1300	rBV	359520	413706	20.36%	3.128%
40	15.883	1310	1312	1314	rVB	69326	84649	4.17%	0.640%
41	16.431	1358	1360	1364	rVB3	57368	94774	4.66%	0.717%
42	17.403	1442	1445	1451	rVB	332113	452424	22.26%	3.421%

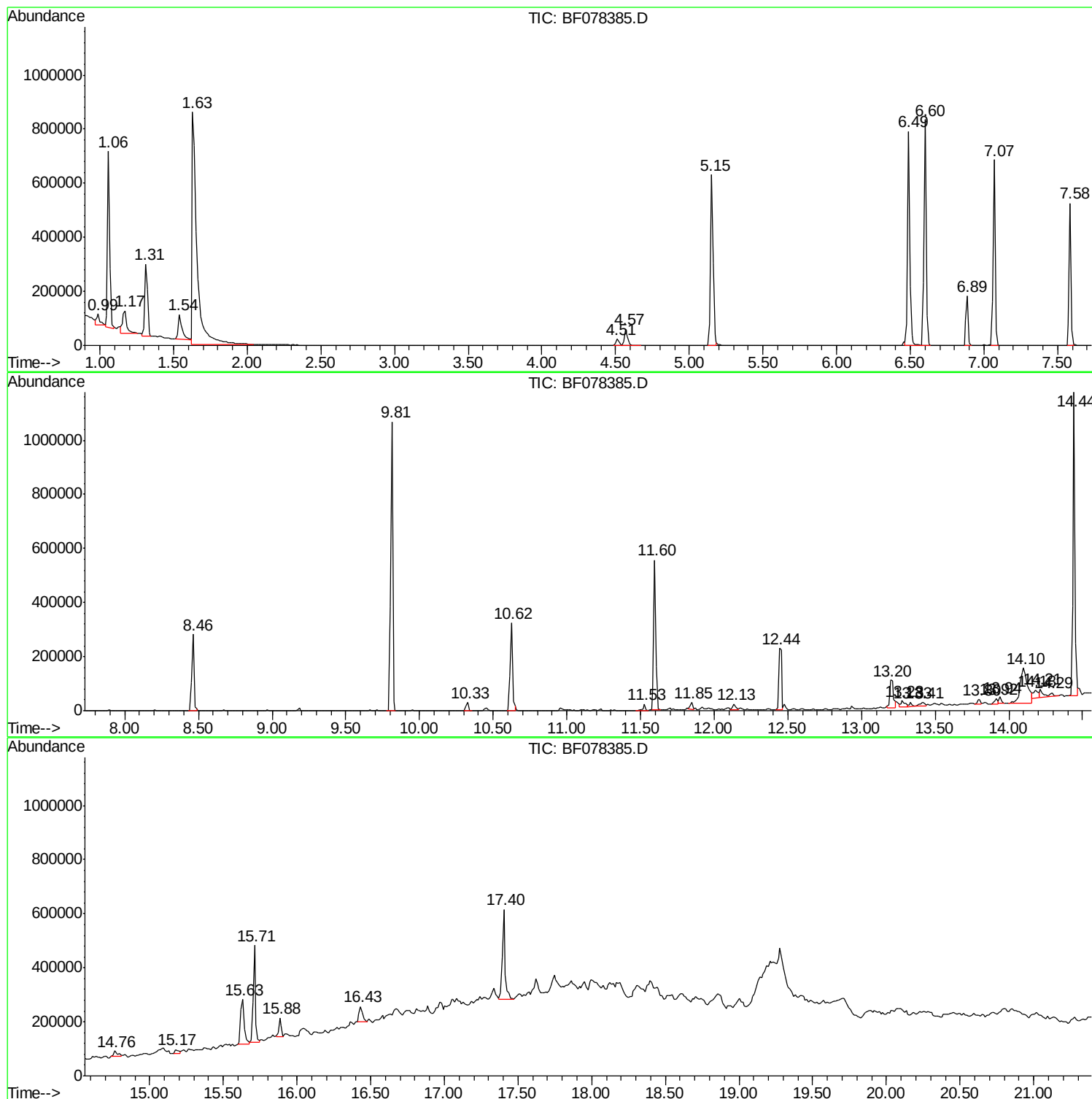
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ClientSampled :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF040115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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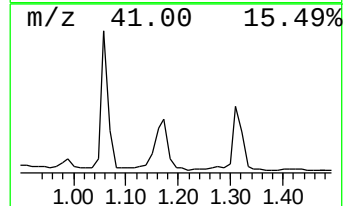
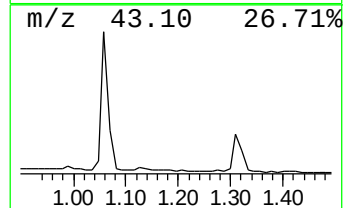
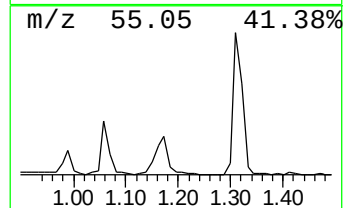
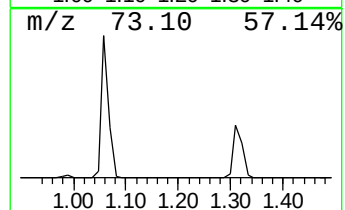
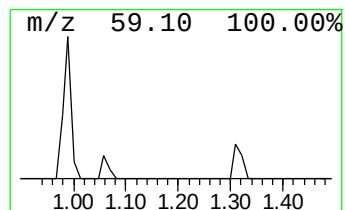
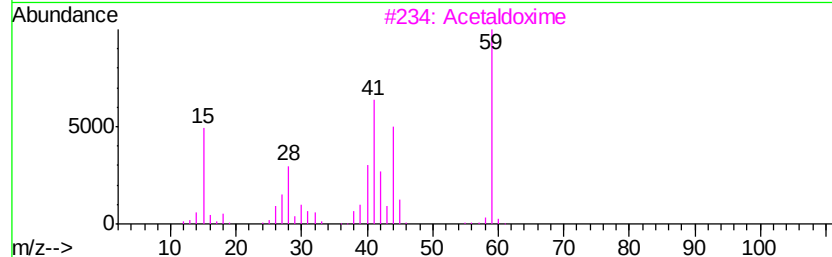
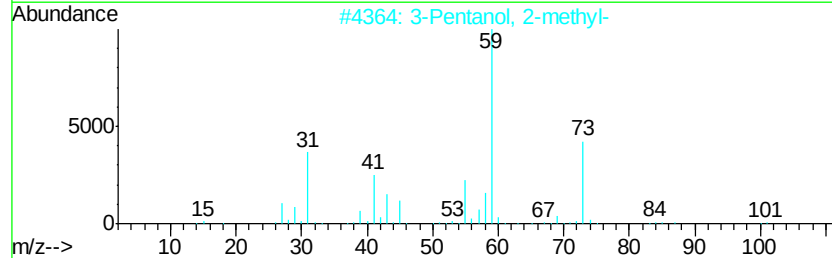
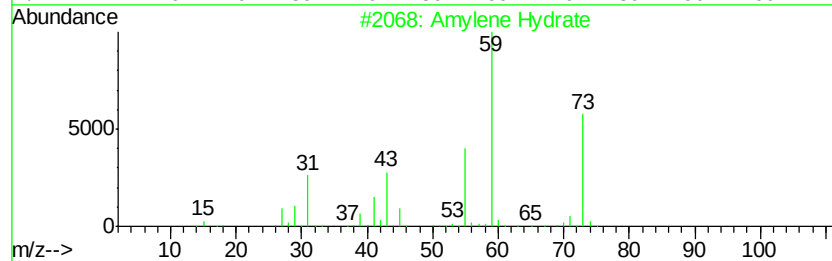
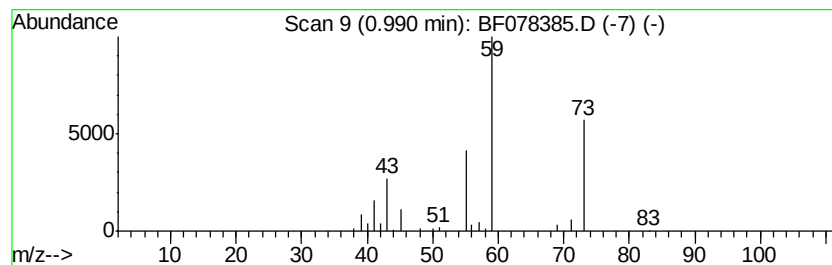
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Amylene Hydrate Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
0.99	5.94 ng	58215	1,4-Dichlorobenzene-d4	6.89

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Amylene Hydrate	88	C5H12O	000075-85-4	78
2		3-Pentanol, 2-methyl-	102	C6H14O	000565-67-3	9
3		Acetaldoxime	59	C2H5NO	000107-29-9	4
4		1,2-Butanediol, (.+/-.)-	90	C4H10O2	026171-83-5	4
5		Butane, 2-methoxy-	88	C5H12O	006795-87-5	4



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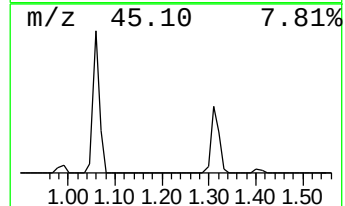
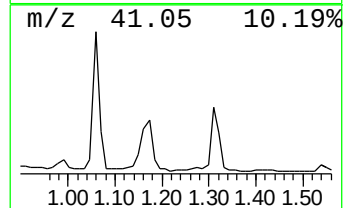
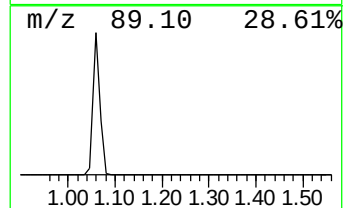
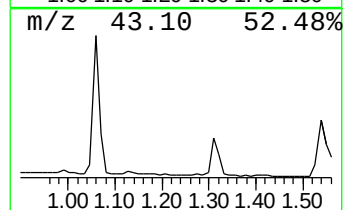
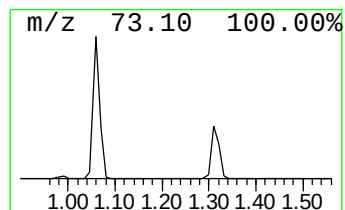
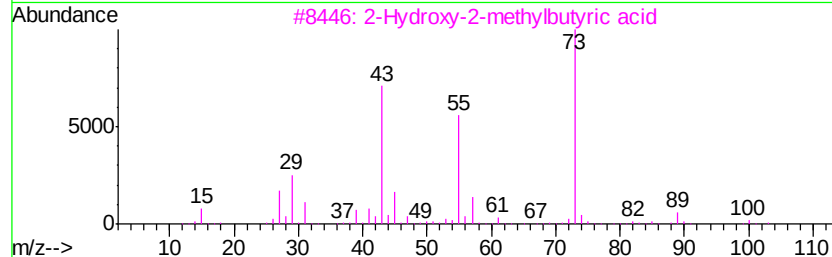
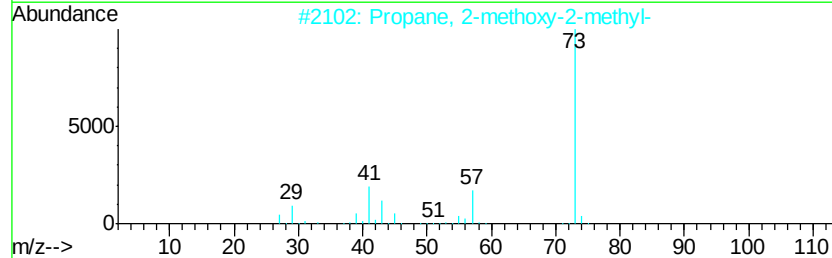
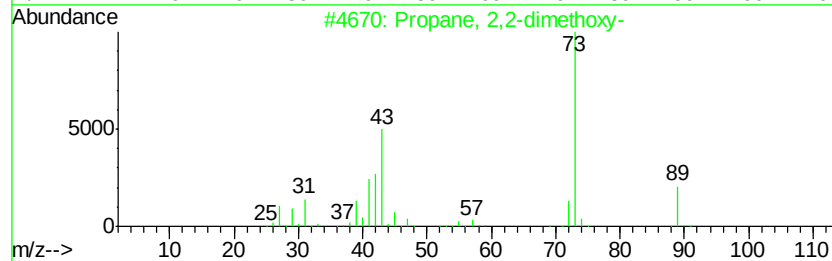
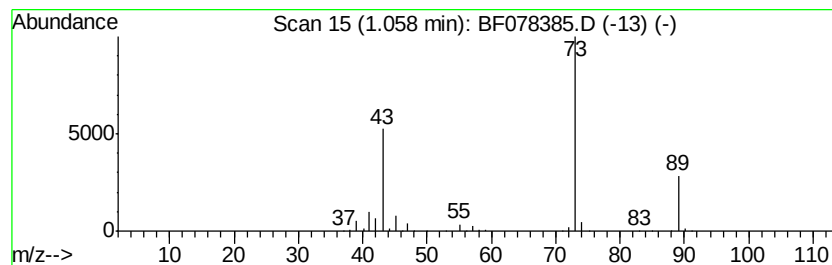
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 2 unknown1.06 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.06	64.61 ng	633566	1,4-Dichlorobenzene-d4	6.89

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Propane, 2,2-dimethoxy-	104	C5H12O2	000077-76-9	38
2		Propane, 2-methoxy-2-methyl-	88	C5H12O	001634-04-4	9
3		2-Hydroxy-2-methylbutyric acid	118	C5H10O3	003739-30-8	9
4		Propanoic acid, 2-methyl-, propy...	130	C7H14O2	000644-49-5	9
5		1-Hexen-4-ol, 1-chloro-3,5-dimet...	162	C8H15ClO	100244-09-5	9



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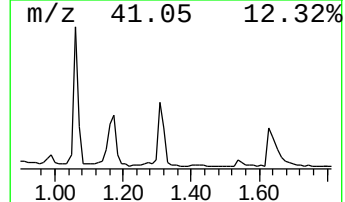
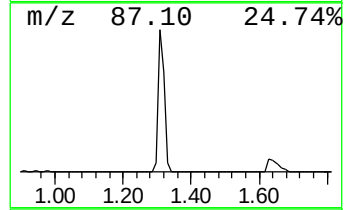
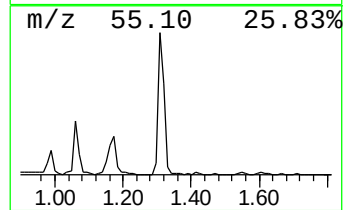
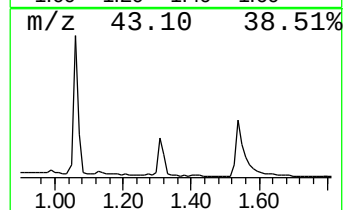
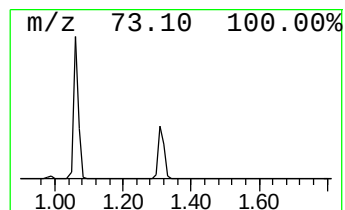
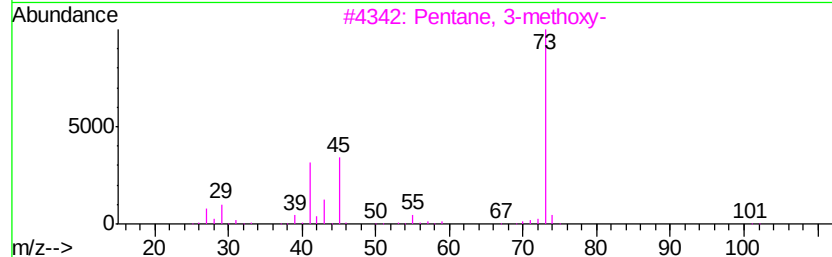
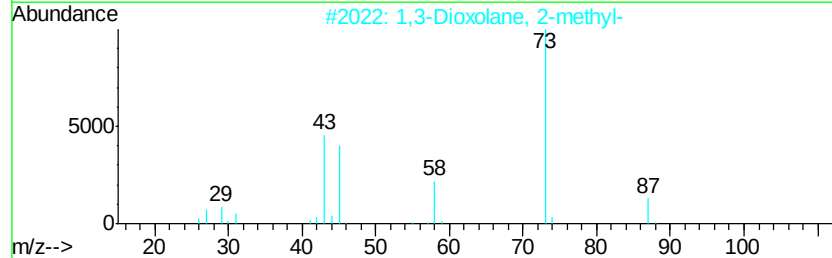
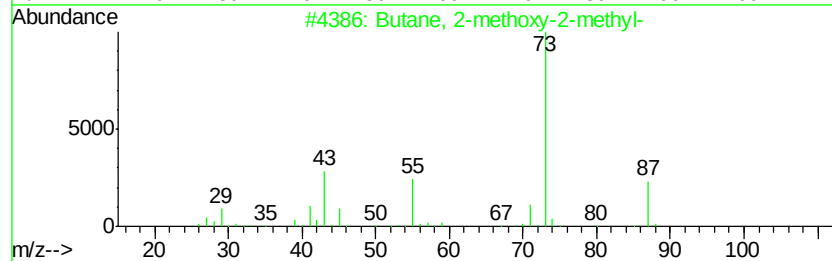
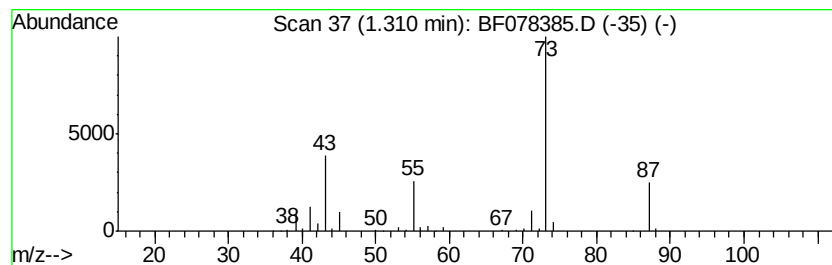
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Butane, 2-methoxy-2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.31	33.11 ng	324697	1,4-Dichlorobenzene-d4	6.89

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2			1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	25
3			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	12
4			Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	12
5			Silane, tetramethyl-	88	C4H12Si	000075-76-3	10



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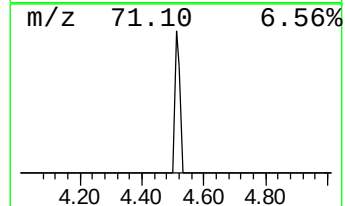
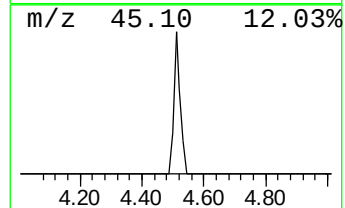
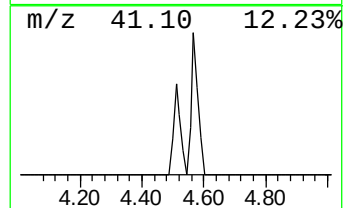
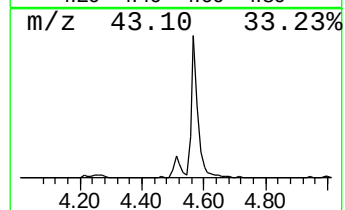
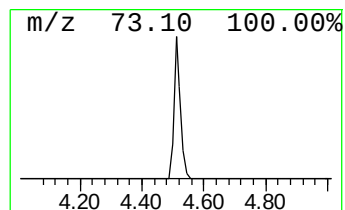
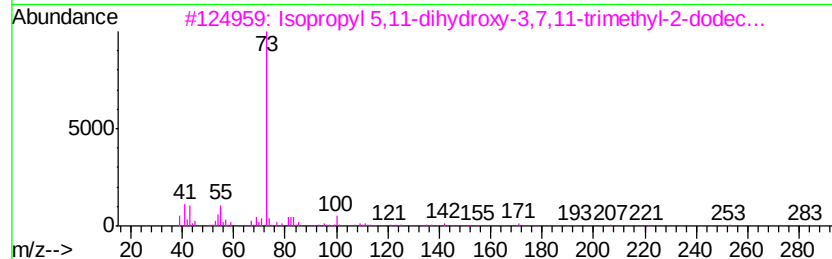
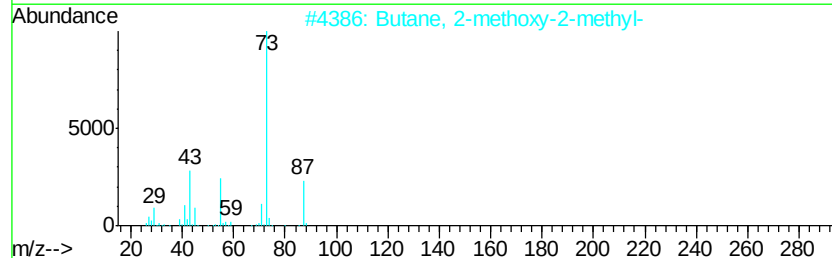
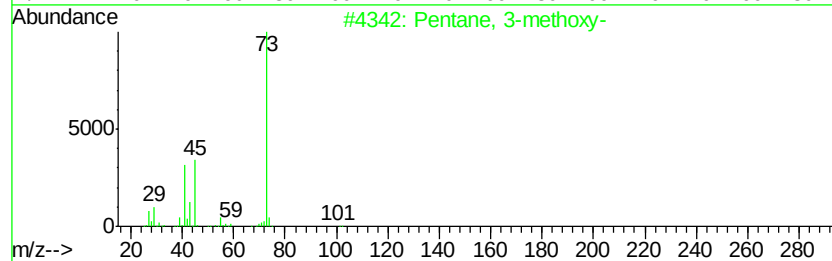
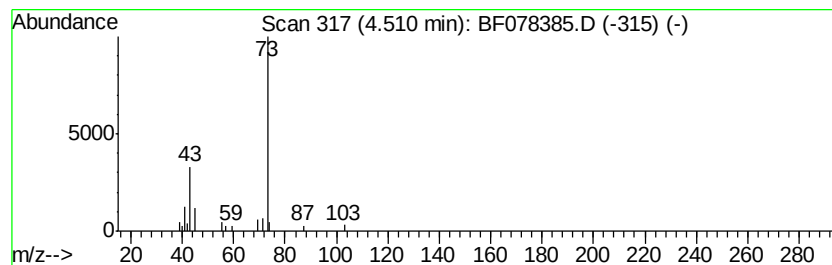
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 unknown4.51 Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.51	3.40 ng	33302	1,4-Dichlorobenzene-d4	6.89

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentane, 3-methoxy-	102	C6H14O	036839-67-5	28
2		Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	9
3		Isopropyl 5,11-dihydroxy-3,7,11-...	314	C18H34O4	1000223-34-1	9
4		Silane, tetramethyl-	88	C4H12Si	000075-76-3	9
5		3-Pentanol, 2,4-dimethyl-	116	C7H16O	000600-36-2	9



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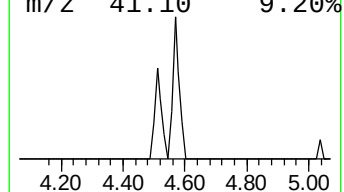
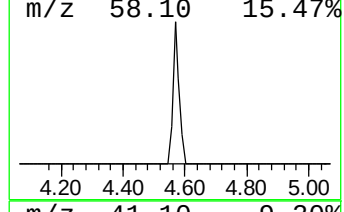
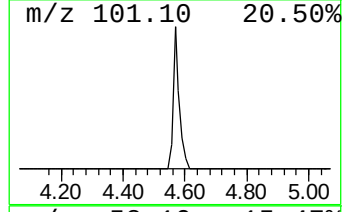
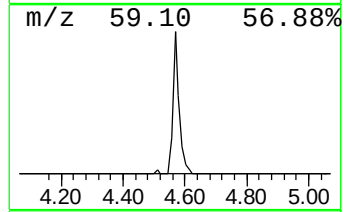
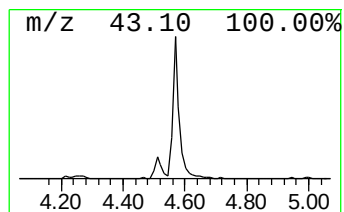
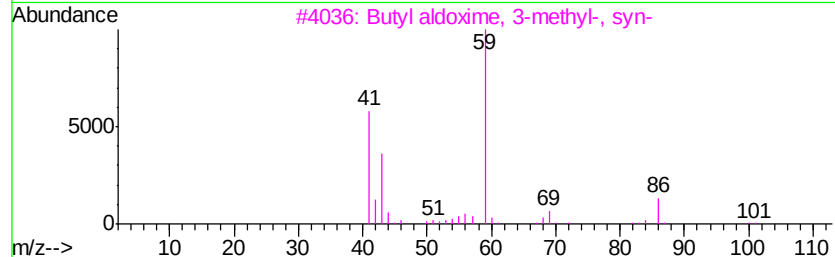
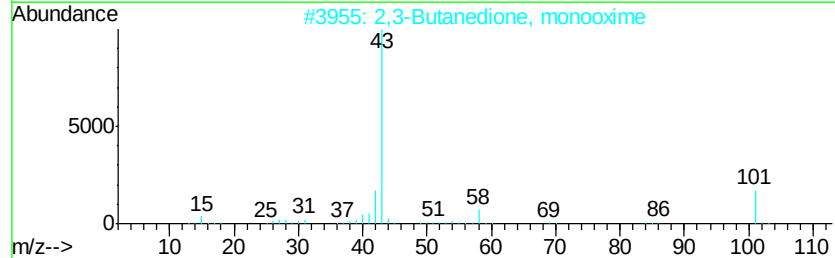
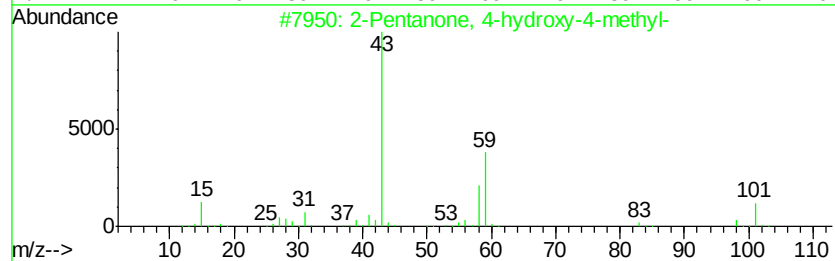
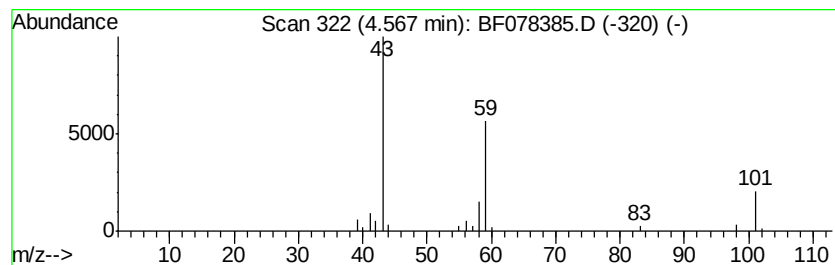
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Peak Number 8 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.57	8.46 ng	82926	1,4-Dichlorobenzene-d4	6.89

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	35
3			Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	17
4			Propylamine	59	C3H9N	000107-10-8	9
5			5-Hexen-2-one	98	C6H10O	000109-49-9	9



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF040915\
Data File : BF078385.D
Acq On : 10 Apr 2015 5:54
Operator : TP/IZ
Sample : G1782-05
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
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ClientSampleId :
SB10(2)

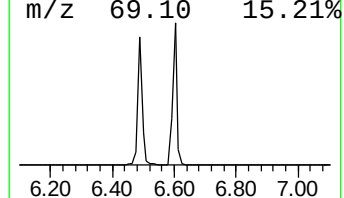
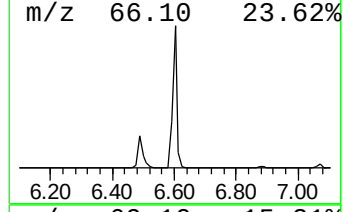
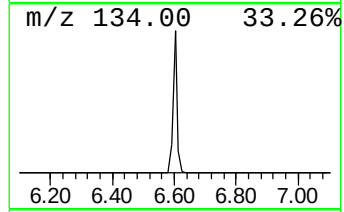
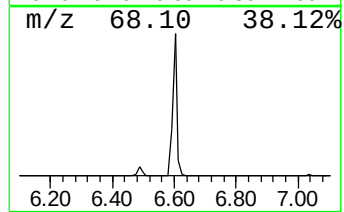
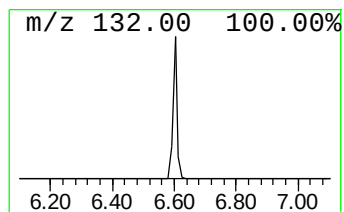
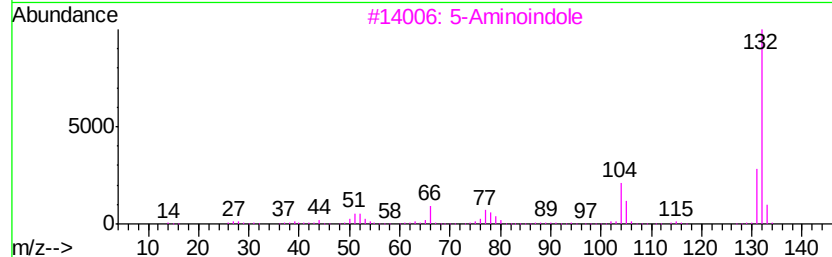
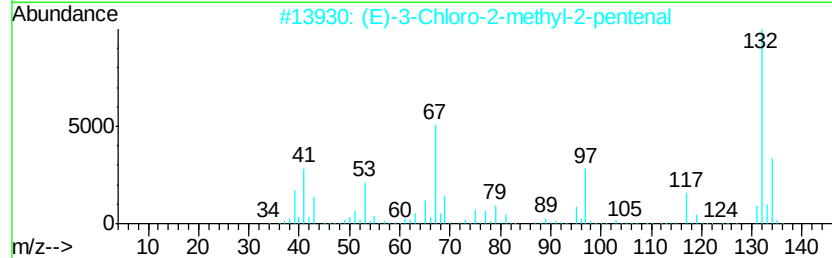
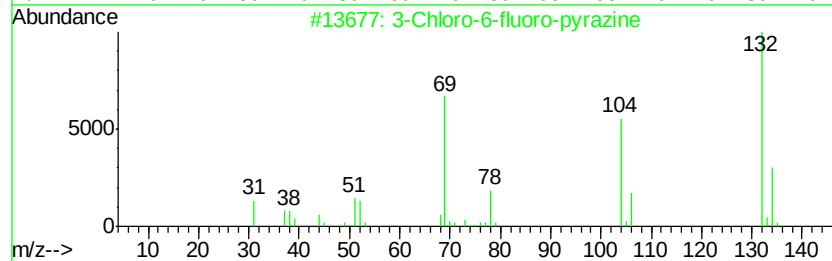
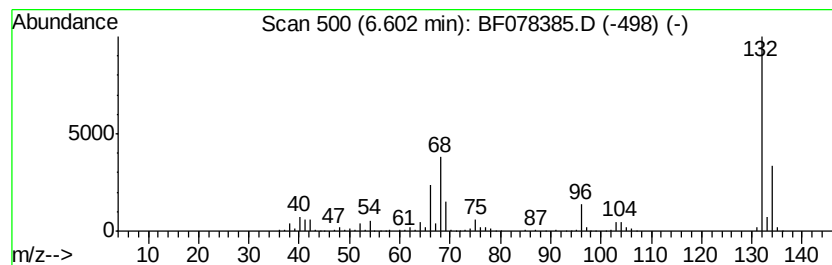
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 unknown6.60 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.60	82.75 ng	811386	1,4-Dichlorobenzene-d4	6.89

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	16
3		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
5		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF040915\
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Acq On : 10 Apr 2015 5:54
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Misc :
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ClientSampleId :
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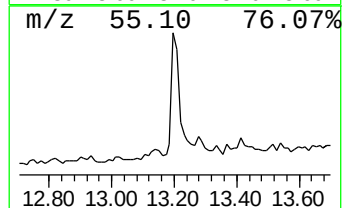
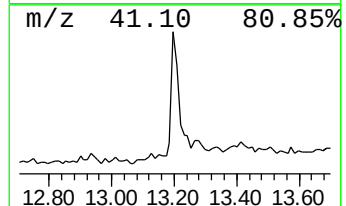
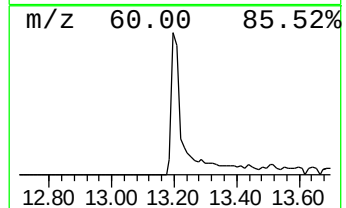
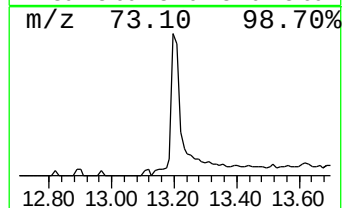
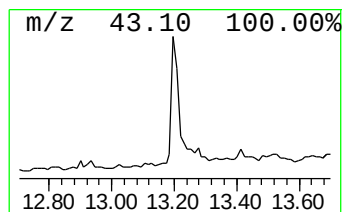
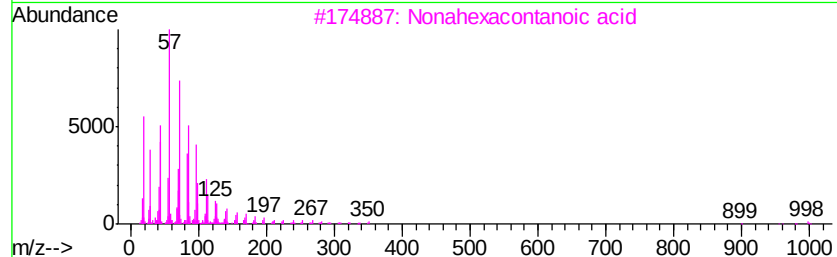
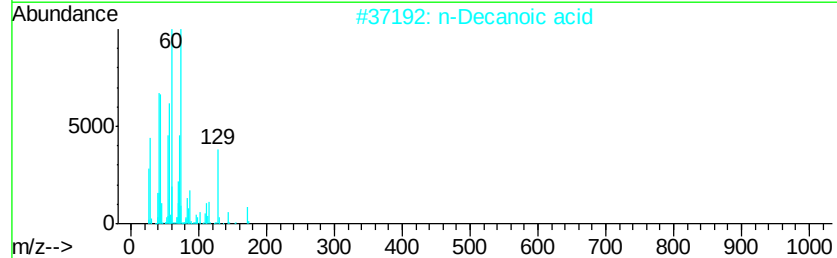
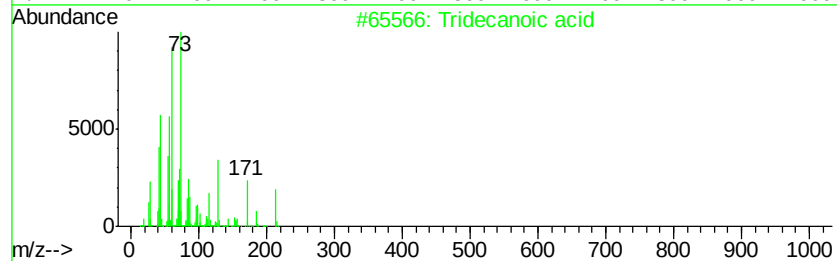
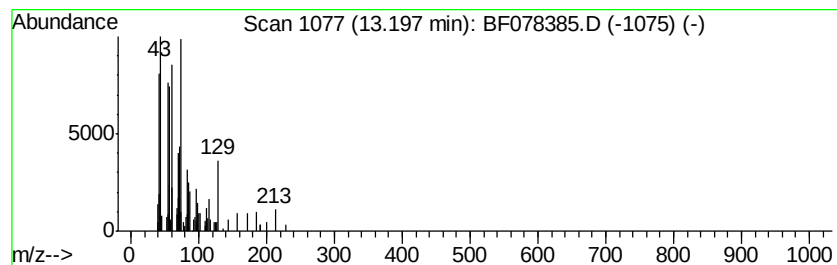
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Tridecanoic acid Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.20	11.54 ng	182809	Phenanthrene-d10	12.45

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tridecanoic acid	214	C13H26O2	000638-53-9	80
2		n-Decanoic acid	172	C10H20O2	000334-48-5	59
3		Nonahexacontanoic acid	999	C69H138O2	040710-32-5	14
4		1-Nonadecanol	284	C19H40O	001454-84-8	11
5		1-Octanol, 2-butyl-	186	C12H26O	003913-02-8	11



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF040915\
Data File : BF078385.D
Acq On : 10 Apr 2015 5:54
Operator : TP/IZ
Sample : G1782-05
Misc :
ALS Vial : 26 Sample Multiplier: 1

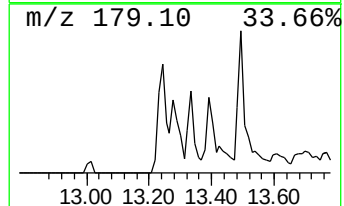
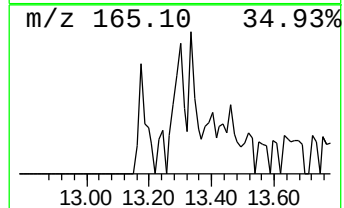
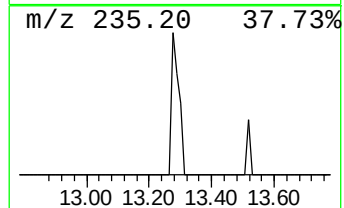
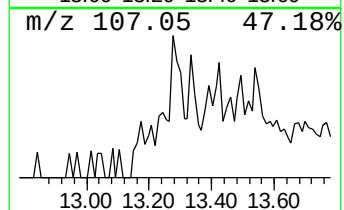
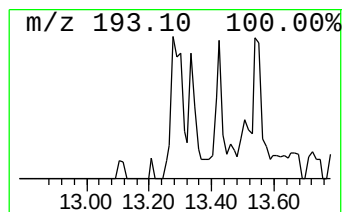
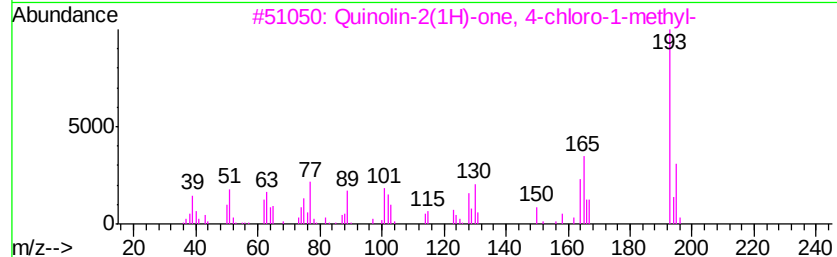
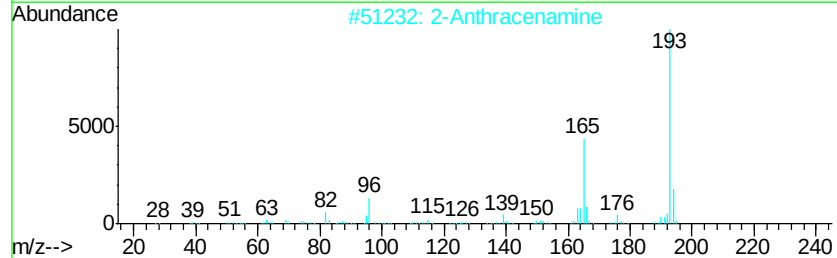
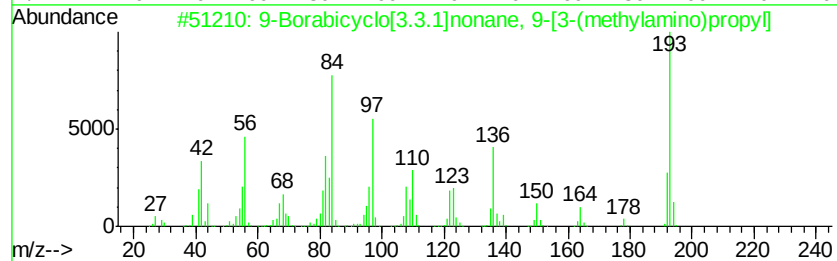
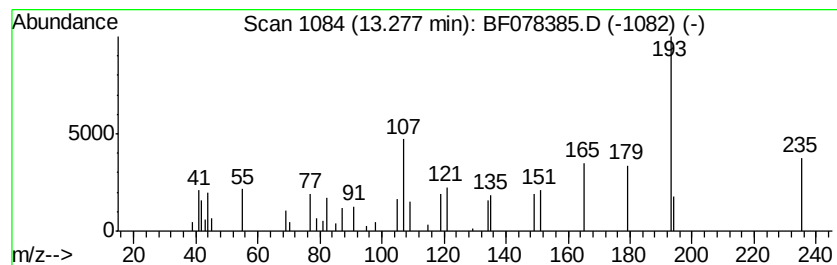
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF040115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 unknown13.28 Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.28	3.05 ng	48338	Phenanthrene-d10	12.45	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9-Borabicyclo[3.3.1]nonane, 9-[3...	193	C12H24BN	1000162-56-0	27
2	2-Anthracenamine	193	C14H11N	000613-13-8	27
3	Quinolin-2(1H)-one, 4-chloro-1-m...	193	C10H8ClNO	032262-17-2	27
4	9-Phenanthrenamine	193	C14H11N	000947-73-9	27
5	1-[1-Methoxy-3,3-dimethyl-2-(3-m...	236	C15H24O2	1000192-05-6	25



Data Path : Z:\HPCHEM1\BNA F\DATA\BF040915\
Data File : BF078385.D
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Operator : TP/IZ
Sample : G1782-05
Misc :
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Instrument :
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ClientSampleId :
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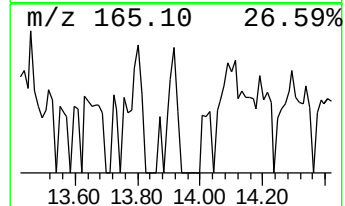
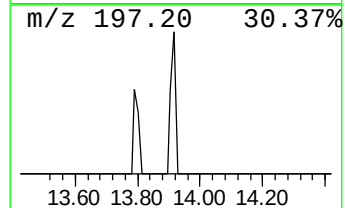
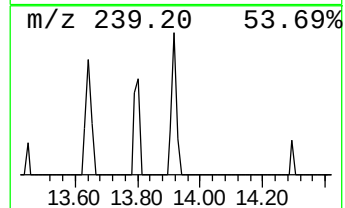
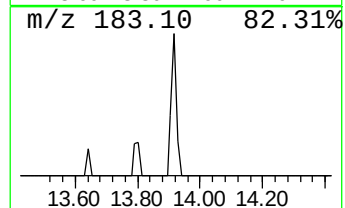
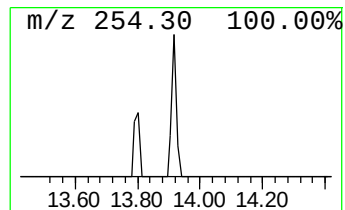
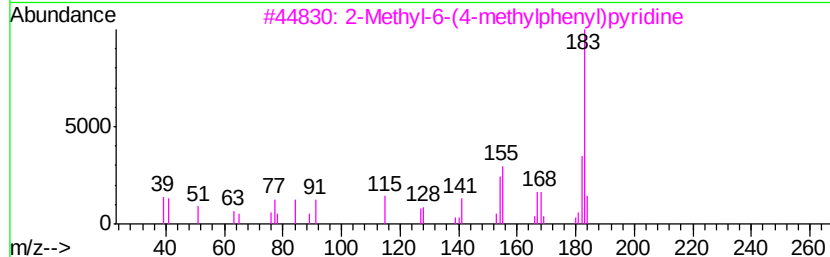
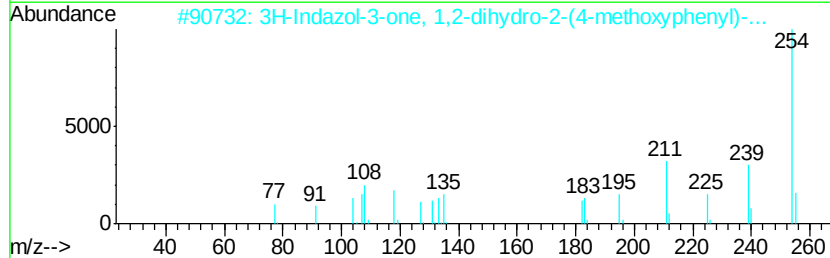
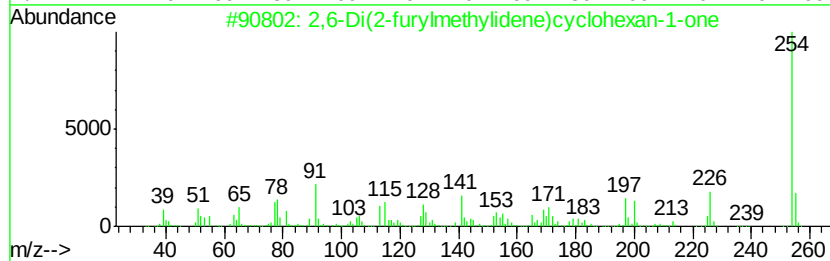
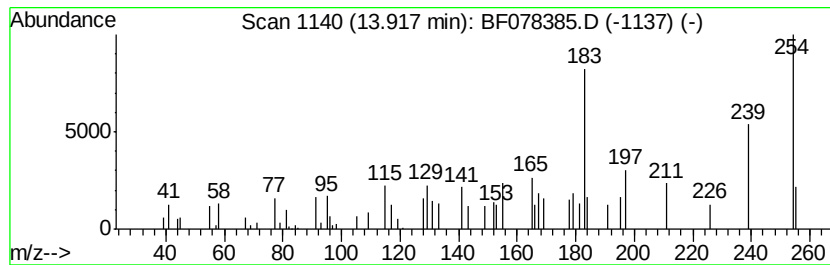
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 unknown13.92 Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.92	2.22 ng	35208	Phenanthrene-d10	12.45

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2,6-Di(2-furvlmethyldene)cyclohexan-1-one	254	C16H14O3	000893-00-5	35
2		3H-Indazol-3-one, 1,2-dihydro-2-(4-methoxyphenyl)-	254	C15H14N2O2	028561-69-5	27
3		2-Methyl-6-(4-methylphenyl)pyridine	183	C13H13N	101893-57-6	22
4		2-Methyl-5-p-tolylpyridine	183	C13H13N	030456-56-5	22
5		Permethrin	390	C21H20Cl2O3	052645-53-1	22



Data Path : Z:\HPCHEM1\BNA F\DATA\BF040915\
Data File : BF078385.D
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Misc :
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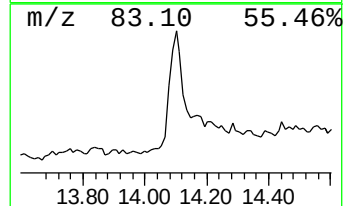
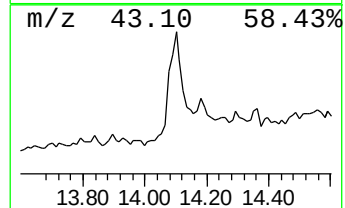
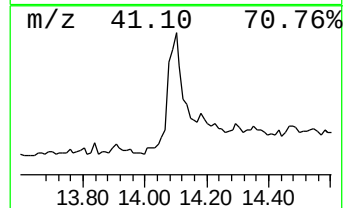
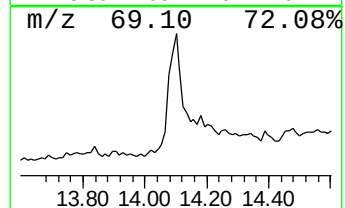
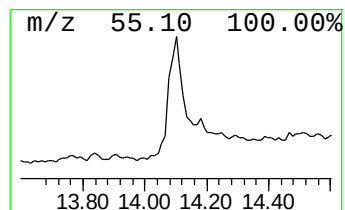
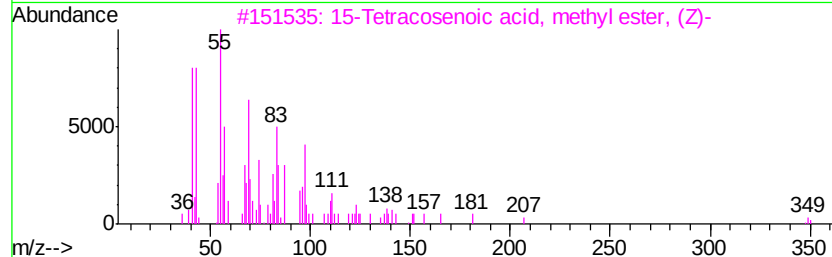
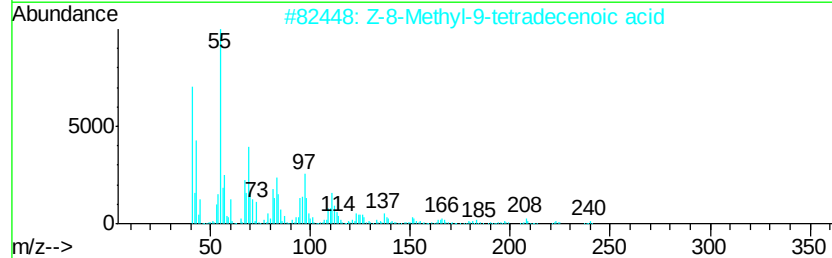
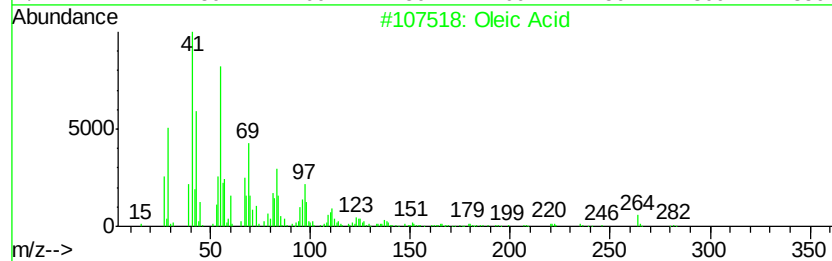
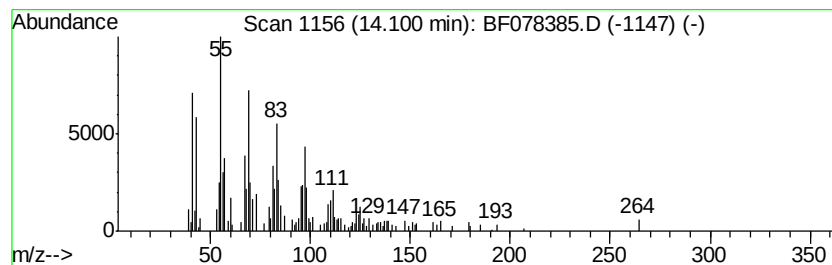
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Oleic Acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.10	23.26 ng	481201	Chrysene-d12	15.71

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Oleic Acid		282	C18H34O2	000112-80-1	91
2	Z-8-Methyl-9-tetradecenoic acid		240	C15H28O2	1000130-84-5	86
3	15-Tetracosenoic acid, methyl es...		380	C25H48O2	002733-88-2	74
4	9-Hexadecenoic acid		254	C16H30O2	002091-29-4	62
5	E-9-Tetradecenoic acid		226	C14H26O2	1000131-35-8	58



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF040915\
Data File : BF078385.D
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Misc :
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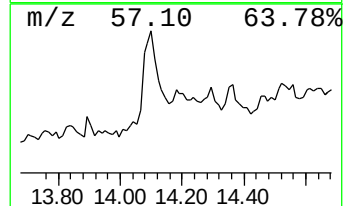
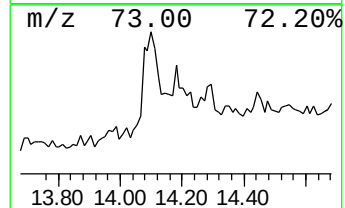
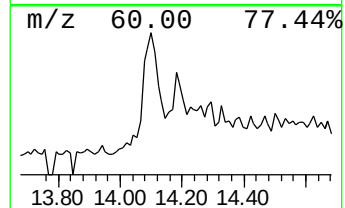
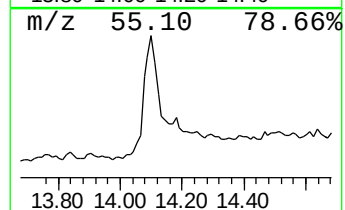
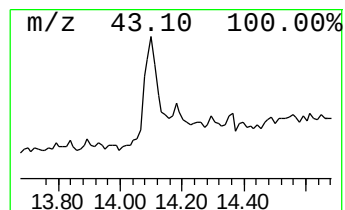
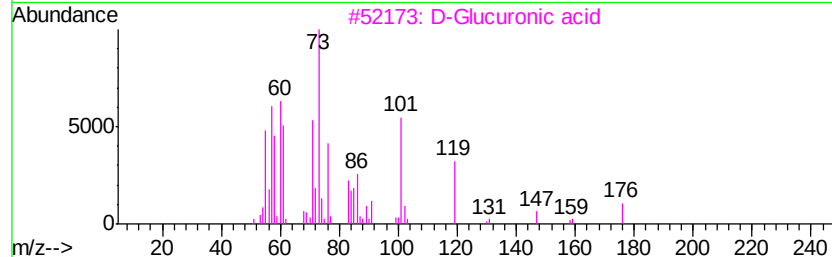
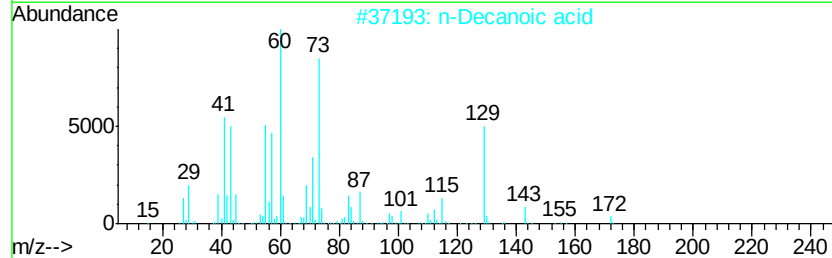
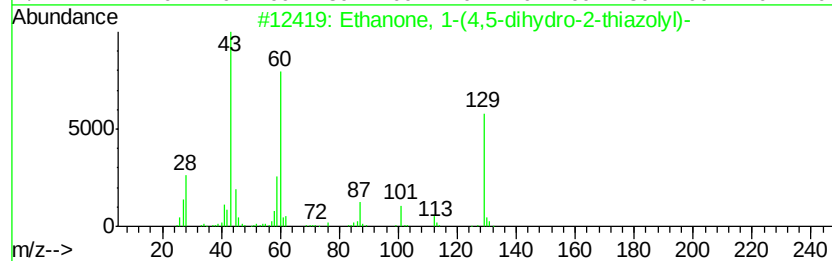
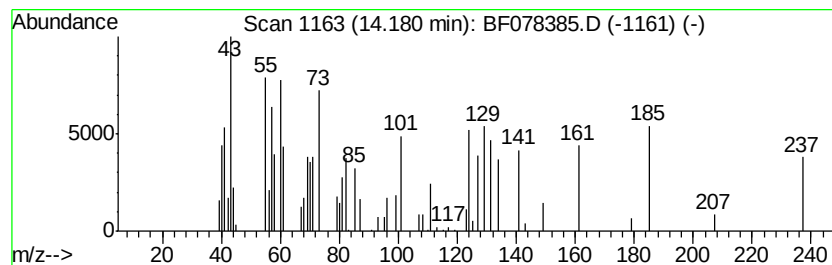
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 16 unknown14.18 Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.18	2.99 ng	61772	Chrysene-d12	15.71

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Ethanone, 1-(4,5-dihydro-2-thiaz...	129	C5H7NOS	029926-41-8	14
2		n-Decanoic acid	172	C10H20O2	000334-48-5	14
3		D-Glucuronic acid	194	C6H10O7	006556-12-3	10
4		D-Mannopyranose	180	C6H12O6	000530-26-7	10
5		Desulphosinigrin	279	C10H17NO6S	005115-81-1	9



Data Path : Z:\HPCHEM1\BNA F\DATA\BF040915\
Data File : BF078385.D
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Sample : G1782-05
Misc :
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ClientSampleId :
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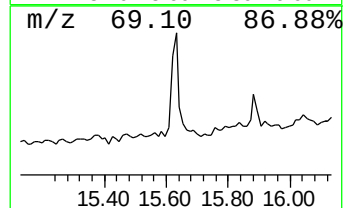
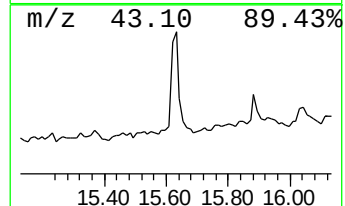
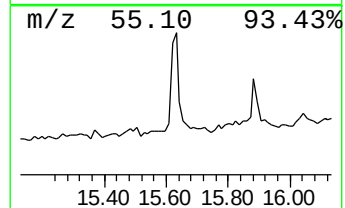
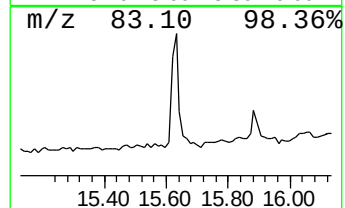
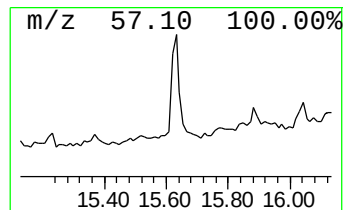
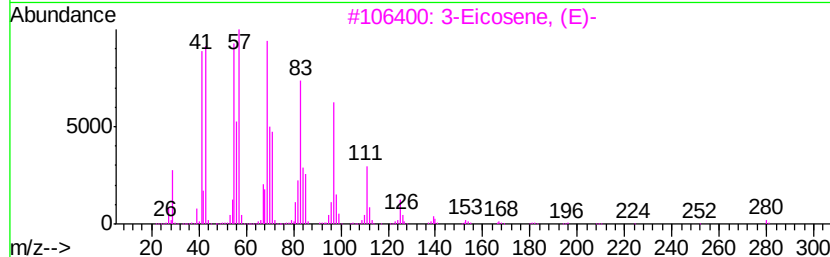
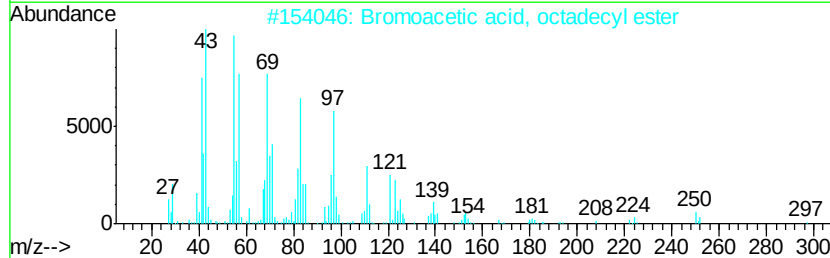
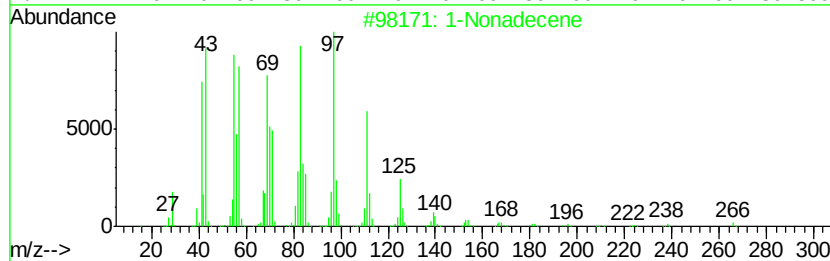
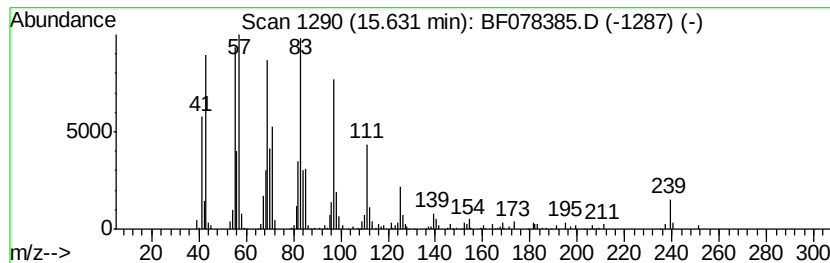
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 1-Nonadecene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.63	13.00 ng	268987	Chrysene-d12	15.71

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Nonadecene	266	C19H38	018435-45-5	91
2		Bromoacetic acid, octadecyl ester	390	C20H39BrO2	018992-03-5	91
3		3-Eicosene, (E)-	280	C20H40	074685-33-9	91
4		Trifluoroacetic acid, n-heptadec...	324	C17H31F3O2	1000216-79-2	91
5		9-Eicosene, (E)-	280	C20H40	074685-29-3	91



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF040915\
Data File : BF078385.D
Acq On : 10 Apr 2015 5:54
Operator : TP/IZ
Sample : G1782-05
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB10(2)

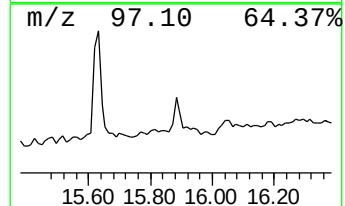
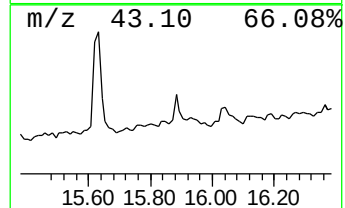
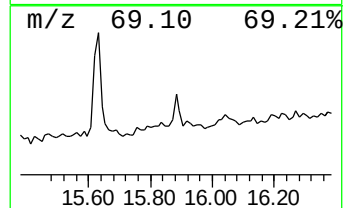
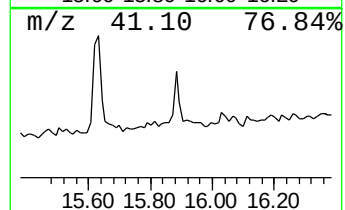
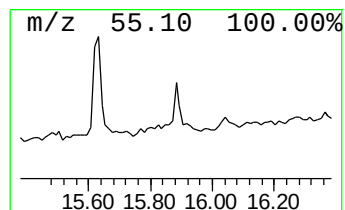
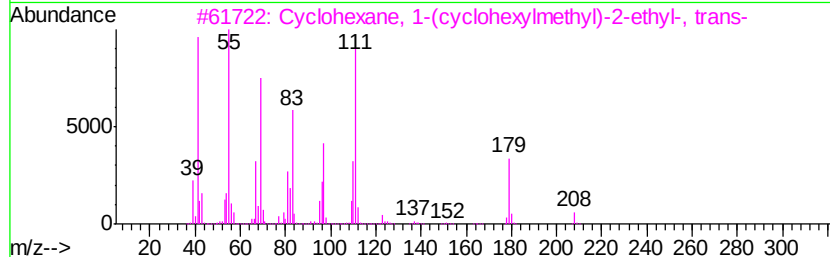
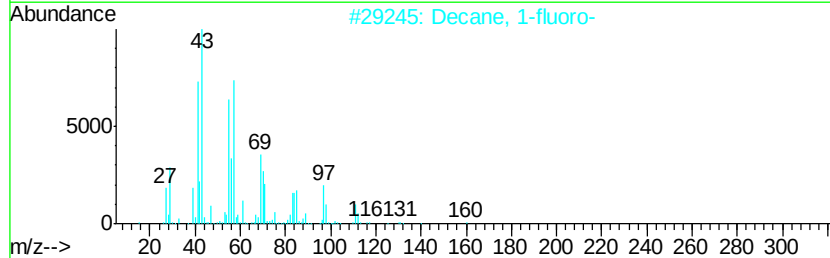
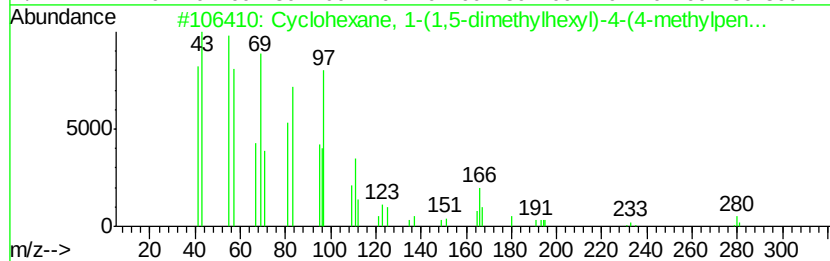
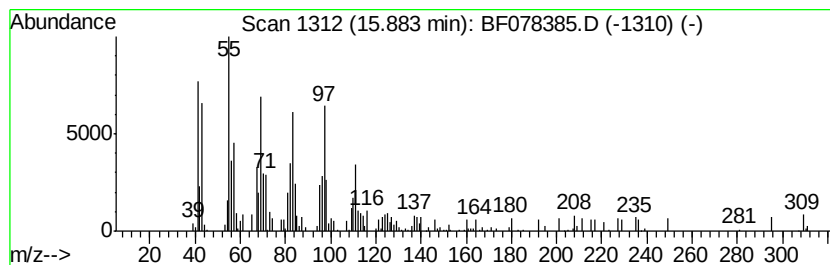
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 19 Cyclohexane, 1-(1,5-dimethy... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.88	4.09 ng	84649	Chrysene-d12	15.71

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, 1-(1,5-dimethylhexyl)-4-(4-methylphenyl)-	280	C20H40	056009-20-2	62
2			Decane, 1-fluoro-	160	C10H21F	000334-56-5	50
3			Cyclohexane, 1-(cyclohexylmethyl)-2-ethyl-, trans-	208	C15H28	054934-92-8	49
4			Cyclopentadecane	210	C15H30	000295-48-7	49
5			Cyclopentadecanone, 2-hydroxy-	240	C15H28O2	004727-18-8	49



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF040915\
Data File : BF078385.D
Acq On : 10 Apr 2015 5:54
Operator : TP/IZ
Sample : G1782-05
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB10(2)

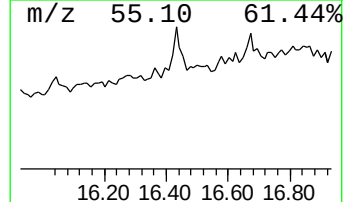
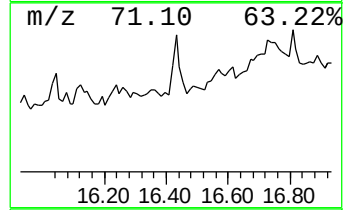
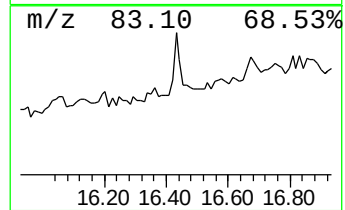
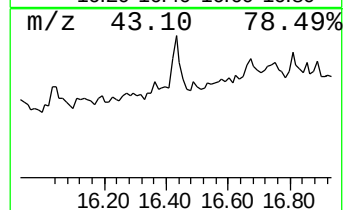
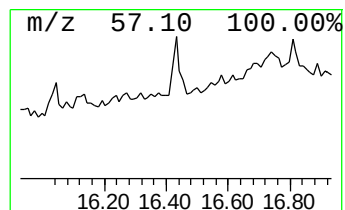
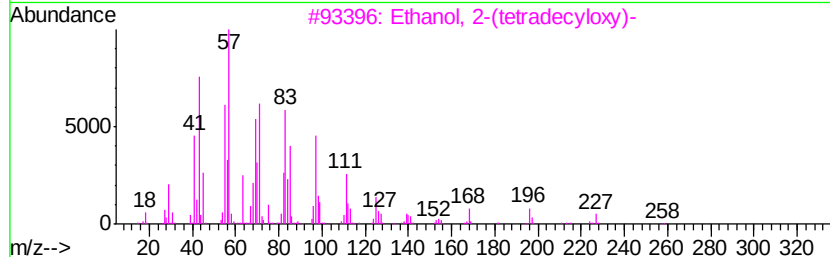
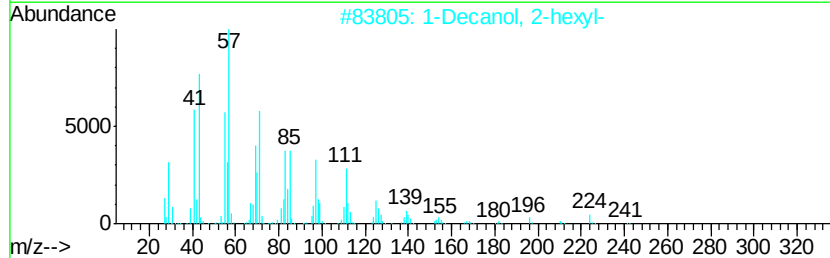
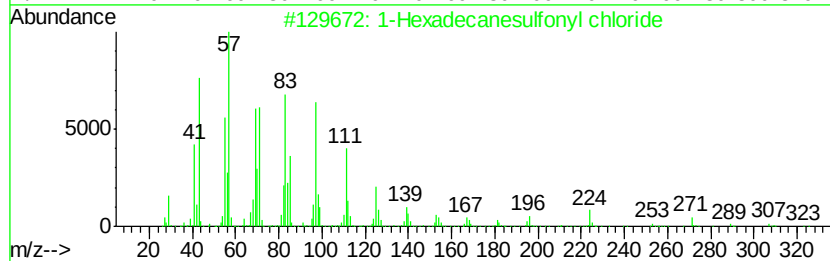
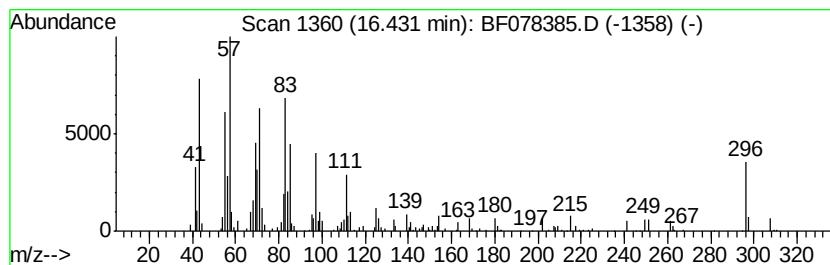
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 20 1-Hexadecanesulfonyl chloride Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.43	4.58 ng	94774	Chrysene-d12	15.71

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Hexadecanesulfonyl chloride	324	C16H33ClO2S	038775-38-1	68
2			1-Decanol, 2-hexyl-	242	C16H34O	002425-77-6	64
3			Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	62
4			1-Heptadecene	238	C17H34	006765-39-5	49
5			Dichloroacetic acid, tetradecyl ...	324	C16H30Cl2O2	083005-02-1	49



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF040915\
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 Sample : G1782-05
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB10(2)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF040115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Amylene Hydrate	0.99	5.9	ng	58215	1	6.89	196116	20.0
unknown1.06	1.06	64.6	ng	633566	1	6.89	196116	20.0
Butane, 2-methoxy...	1.31	33.1	ng	324697	1	6.89	196116	20.0
unknown4.51	4.51	3.4	ng	33302	1	6.89	196116	20.0
2-Pentanone, 4-hy...	4.57	8.5	ng	82926	1	6.89	196116	20.0
unknown6.60	6.60	82.8	ng	811386	1	6.89	196116	20.0
Tridecanoic acid	13.20	11.5	ng	182809	4	12.45	316803	20.0
unknown13.28	13.28	3.0	ng	48338	4	12.45	316803	20.0
unknown13.92	13.92	2.2	ng	35208	4	12.45	316803	20.0
Oleic Acid	14.10	23.3	ng	481201	5	15.71	413706	20.0
unknown14.18	14.18	3.0	ng	61772	5	15.71	413706	20.0
1-Nonadecene	15.63	13.0	ng	268987	5	15.71	413706	20.0
Cyclohexane, 1-(1...	15.88	4.1	ng	84649	5	15.71	413706	20.0
1-Hexadecanesulfo...	16.43	4.6	ng	94774	5	15.71	413706	20.0