

Data Path : Z:\HPCHEM1\BNA F\DATA\BF101515\
 Data File : BF082288.D
 Acq On : 14 Oct 2015 19:39
 Operator : UM/IZ
 Sample : G4015-11
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 LF-7

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_F\METHODS\8270-BF101015.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	4.309	187	190	201	rBV	58698	86529	3.70%	0.516%
2	4.983	246	249	252	rBV	745720	936477	39.99%	5.587%
3	5.406	284	286	289	rBV	1087644	1039853	44.41%	6.203%
4	6.457	375	378	380	rBV	1347690	1381014	58.98%	8.239%
5	6.583	387	389	392	rBV	1306492	1225367	52.33%	7.310%
6	6.823	407	410	412	rBV	364546	358037	15.29%	2.136%
7	6.972	421	423	426	rBV	1003073	1052811	44.96%	6.281%
8	7.395	457	460	462	rBV	738994	854457	36.49%	5.097%
9	8.115	520	523	525	rBV	371977	496500	21.20%	2.962%
10	9.189	614	617	620	rBV	2167820	1941779	82.93%	11.584%
11	9.589	649	652	654	rBV	404968	353837	15.11%	2.111%
12	9.795	667	670	674	rBV	22596	30322	1.29%	0.181%
13	9.863	674	676	679	rBV	660969	665030	28.40%	3.967%
14	10.298	712	714	716	rVB	31795	27567	1.18%	0.164%
15	10.652	743	745	748	rBV2	541882	663535	28.34%	3.958%
16	10.766	753	755	760	rVB	45474	60872	2.60%	0.363%
17	11.224	793	795	796	rBV	32607	39152	1.67%	0.234%
18	11.361	805	807	809	rBV	543402	665636	28.43%	3.971%
19	11.646	830	832	834	rBV	49092	39566	1.69%	0.236%
20	12.058	866	868	870	rBV	29904	27713	1.18%	0.165%
21	12.447	900	902	904	rVB	46979	37674	1.61%	0.225%
22	12.824	933	935	938	rBV2	54805	67817	2.90%	0.405%
23	12.949	944	946	949	rBV	1791866	2341565	100.00%	13.969%
24	13.178	964	966	968	rBV	46654	66710	2.85%	0.398%
25	13.224	968	970	974	rBV2	15220	29814	1.27%	0.178%
26	13.315	976	978	981	rVV	20326	32222	1.38%	0.192%
27	13.407	983	986	988	rBV3	13090	30371	1.30%	0.181%
28	13.544	995	998	1000	rBV	81971	98617	4.21%	0.588%
29	13.887	1025	1028	1036	rVB	274986	422466	18.04%	2.520%
30	14.001	1036	1038	1040	rBV2	14434	29266	1.25%	0.175%
31	14.047	1040	1042	1045	rBV	490905	658090	28.10%	3.926%
32	14.241	1056	1059	1060	rBV	61049	81000	3.46%	0.483%
33	14.344	1066	1068	1072	rBV3	18553	36757	1.57%	0.219%
34	14.584	1086	1089	1094	rBV	65823	111673	4.77%	0.666%

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Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF101015.M
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35	14.687	1096	1098	1104	rVB2	15259	25587	1.09%	0.153%
36	14.927	1114	1119	1121	rBV2	49526	97060	4.15%	0.579%
37	14.995	1123	1125	1127	rBV	27168	37563	1.60%	0.224%
38	15.270	1147	1149	1151	rBV	35133	47726	2.04%	0.285%
39	15.601	1175	1178	1180	rBV	339434	487449	20.82%	2.908%
40	16.070	1216	1219	1221	rBV	16787	28045	1.20%	0.167%
41	16.127	1221	1224	1228	rVB2	25108	49192	2.10%	0.293%

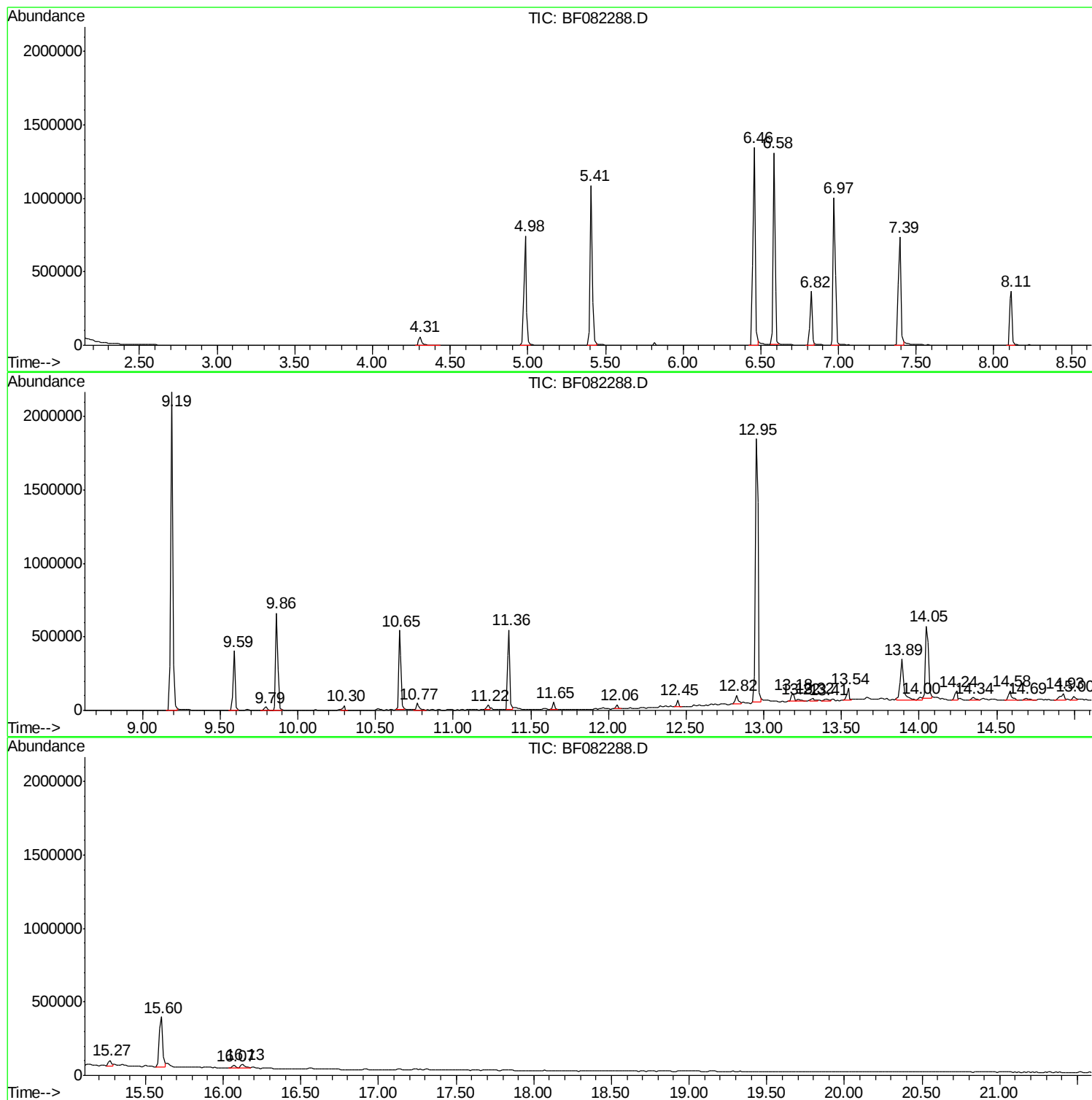
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF101015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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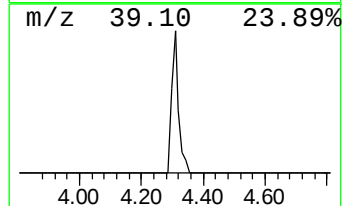
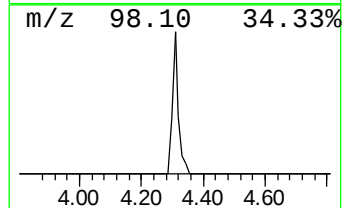
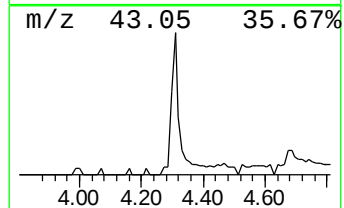
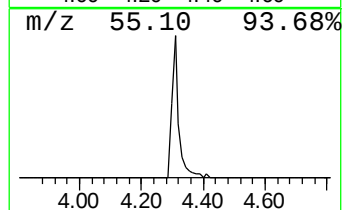
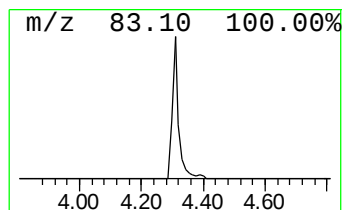
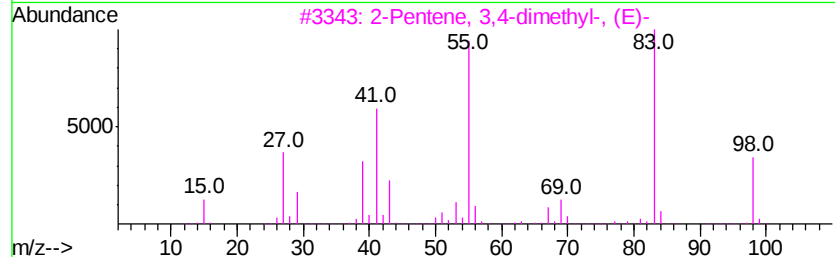
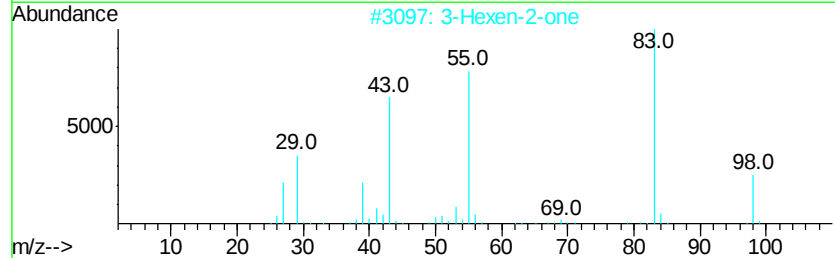
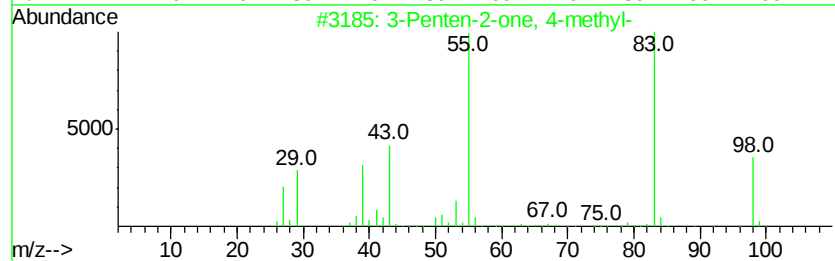
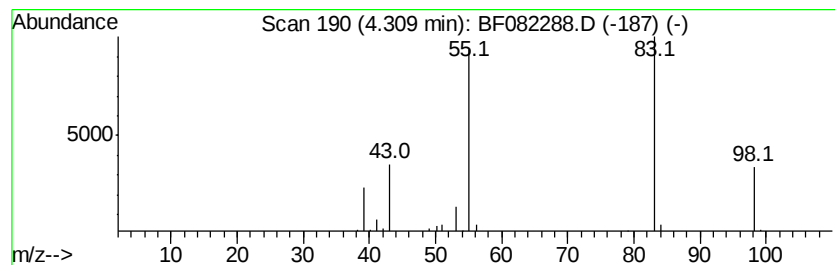
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 3-Penten-2-one, 4-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.31	4.83 ng	86529	1,4-Dichlorobenzene-d4	6.82

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Penten-2-one, 4-methyl-	98	C6H10O	000141-79-7	91
2		3-Hexen-2-one	98	C6H10O	000763-93-9	90
3		2-Pentene, 3,4-dimethyl-, (E)-	98	C7H14	004914-92-5	90
4		2-Pentene, 2,4-dimethyl-	98	C7H14	000625-65-0	86
5		2-Pentene, 3,4-dimethyl-, (Z)-	98	C7H14	004914-91-4	78



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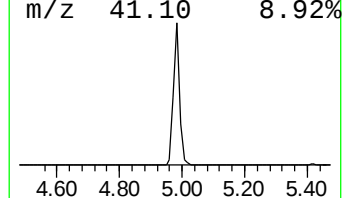
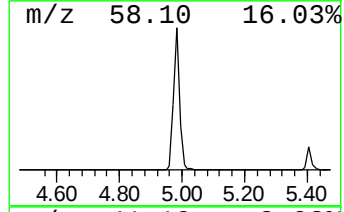
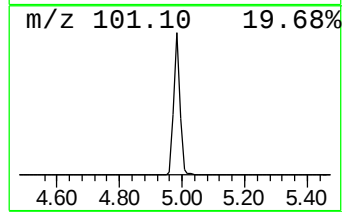
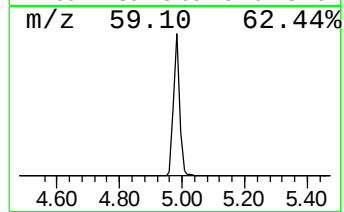
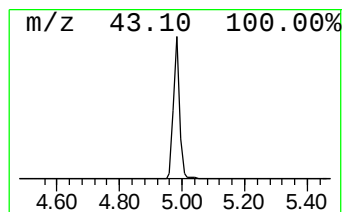
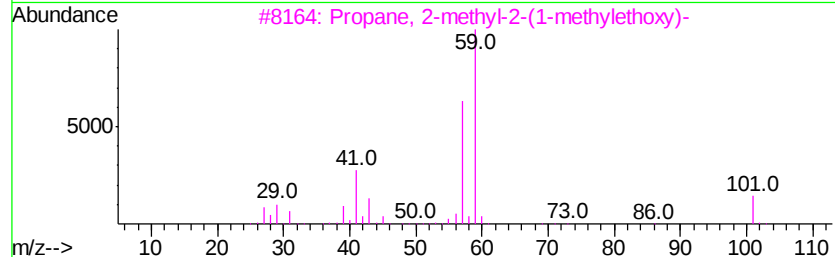
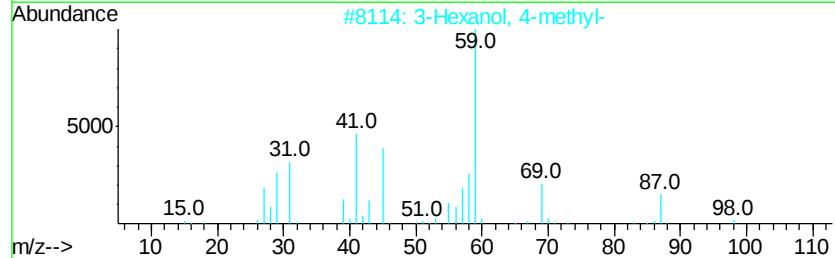
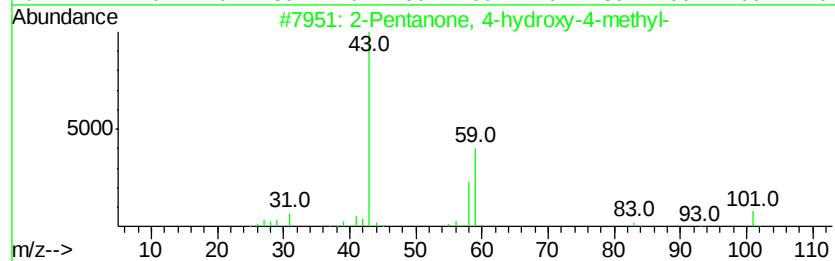
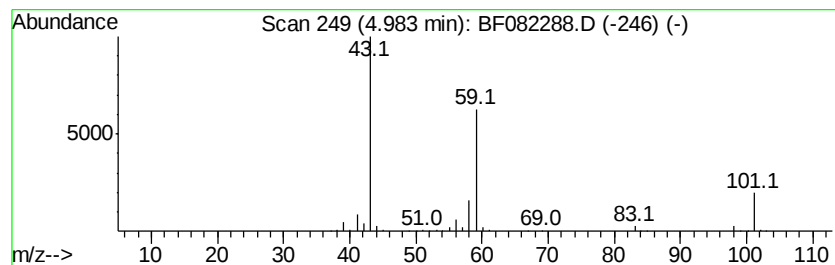
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.98	52.31 ng	936477	1,4-Dichlorobenzene-d4	6.82

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	28
3			Propane, 2-methyl-2-(1-methylethoxy)-	116	C7H16O	017348-59-3	28
4			Acetic acid, cyano-, 1,1-dimethyl-	141	C7H11NO2	001116-98-9	17
5			Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9



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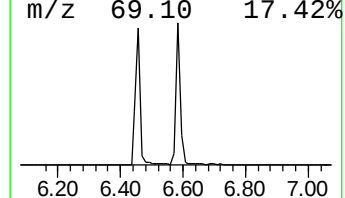
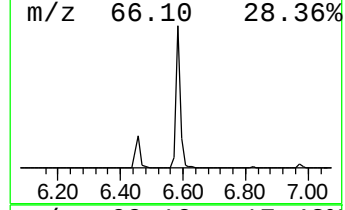
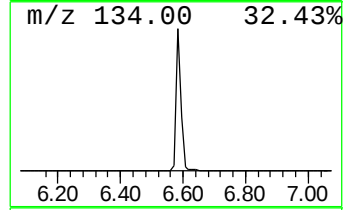
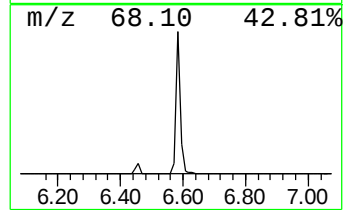
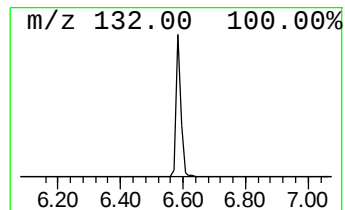
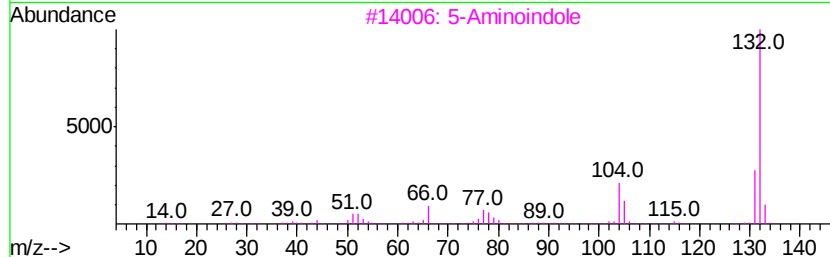
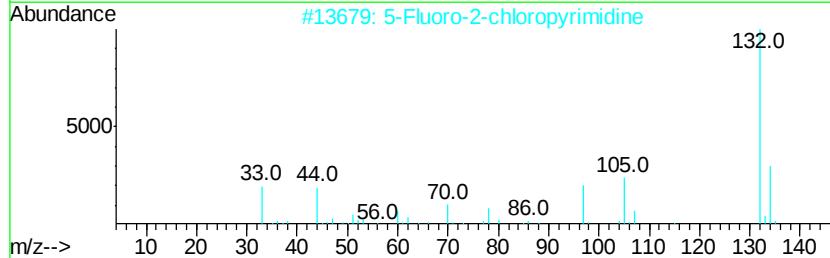
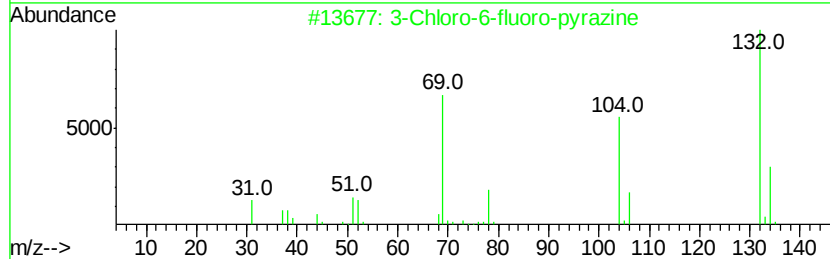
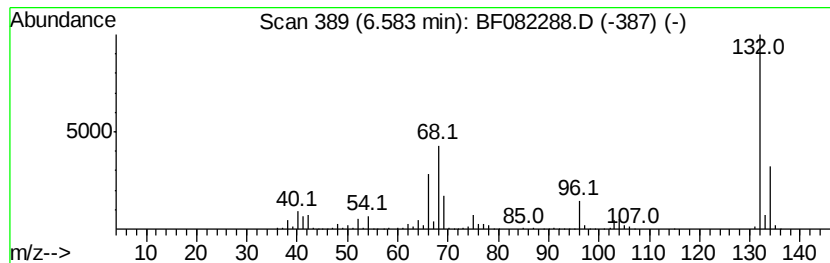
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 3 unknown6.58 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.58	68.45 ng	1225370	1,4-Dichlorobenzene-d4	6.82

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	35
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
3		5-Aminoindole	132	C8H8N2	005192-03-0	12
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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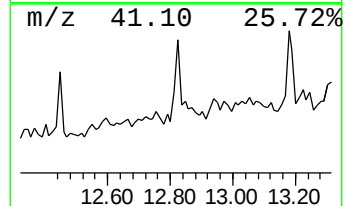
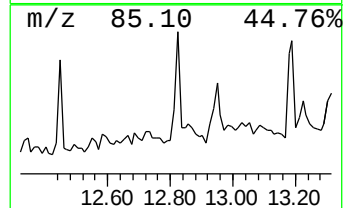
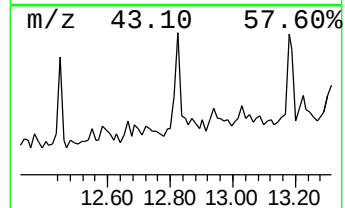
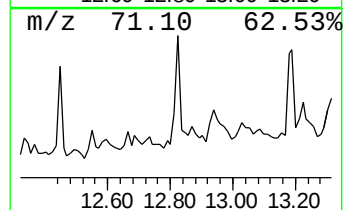
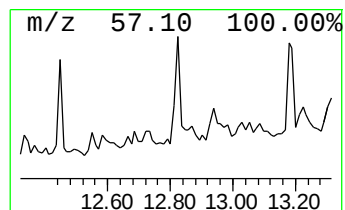
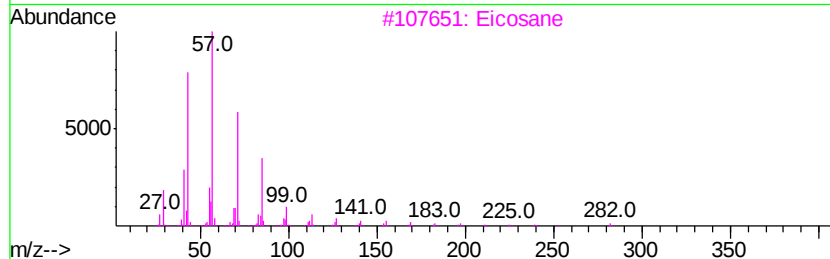
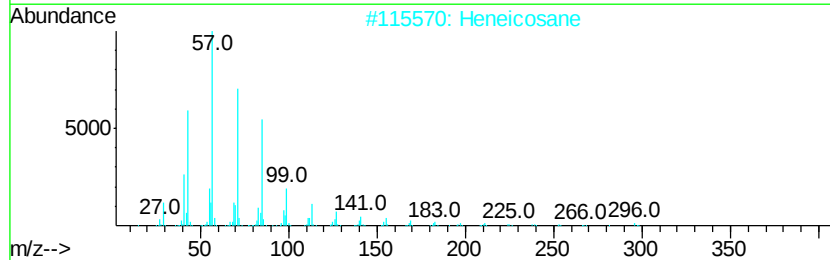
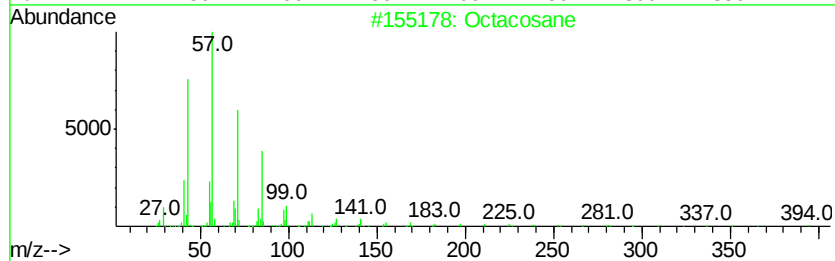
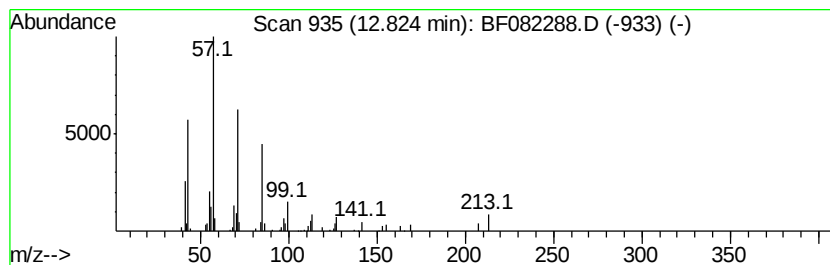
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Octacosane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.82	2.06 ng	67817	Chrysene-d12	14.05

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octacosane	394	C28H58	000630-02-4	87
2		Heneicosane	296	C21H44	000629-94-7	87
3		Eicosane	282	C20H42	000112-95-8	86
4		Pentadecane, 8-heptyl-	310	C22H46	071005-15-7	86
5		Heptacosane	380	C27H56	000593-49-7	86



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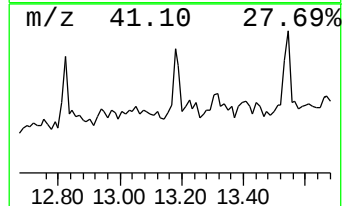
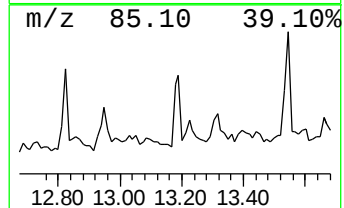
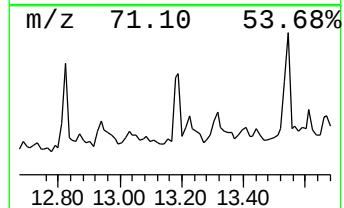
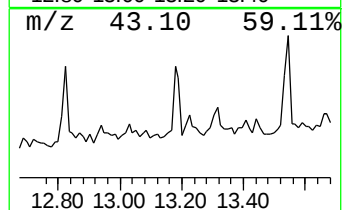
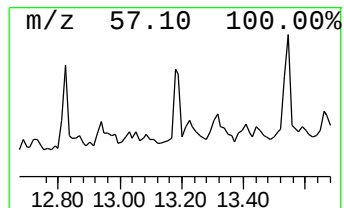
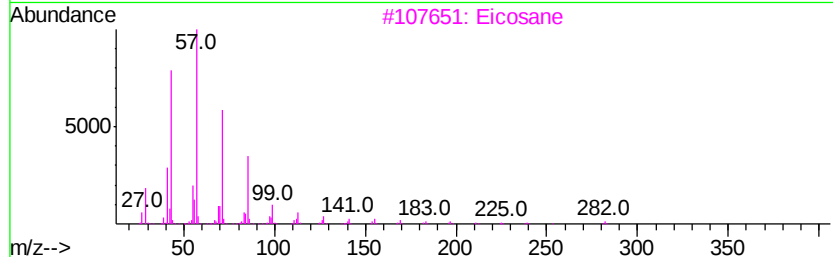
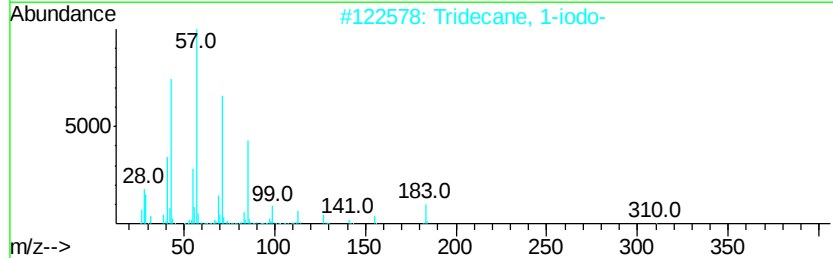
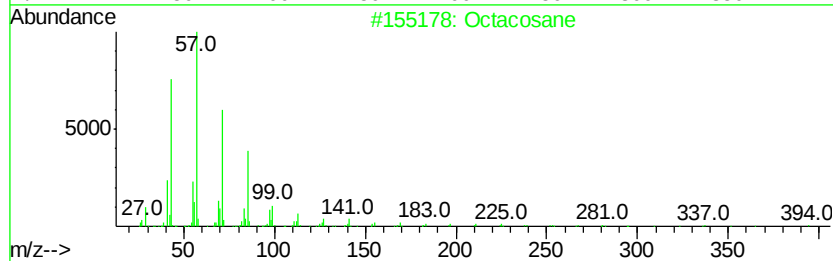
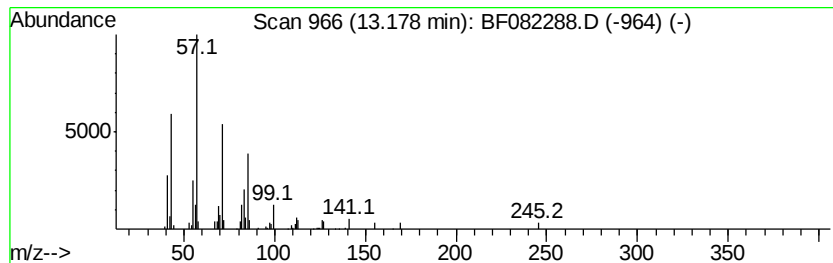
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Tridecane, 1-iodo- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.18	2.03 ng	66710	Chrysene-d12	14.05
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Octacosane	394 C28H58	000630-02-4 83
2		Tridecane, 1-iodo-	310 C13H27I	035599-77-0 74
3		Eicosane	282 C20H42	000112-95-8 72
4		Tetradecane	198 C14H30	000629-59-4 72
5		Dodecane, 2-methyl-6-propyl-	226 C16H34	055045-08-4 72



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF101515\
Data File : BF082288.D
Acq On : 14 Oct 2015 19:39
Operator : UM/IZ
Sample : G4015-11
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
LF-7

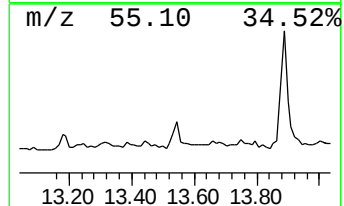
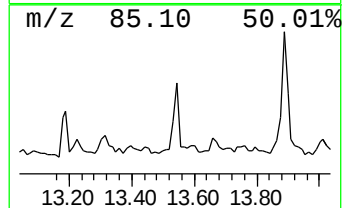
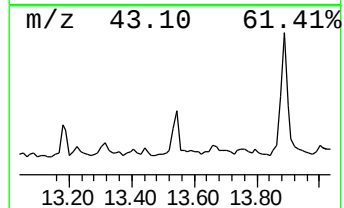
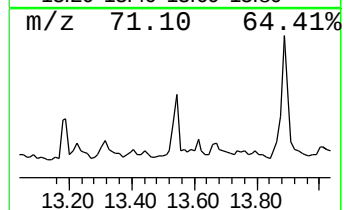
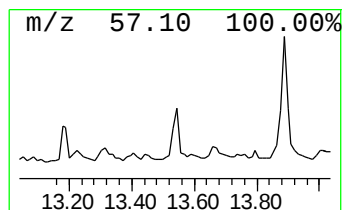
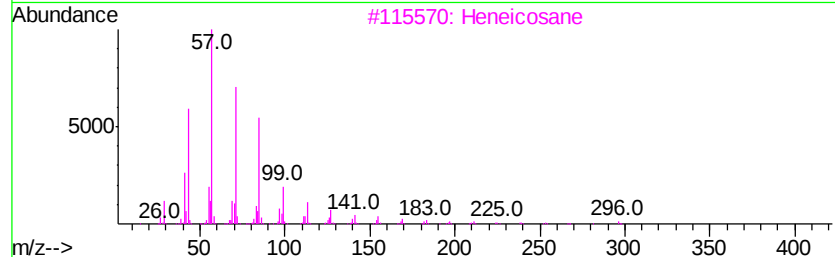
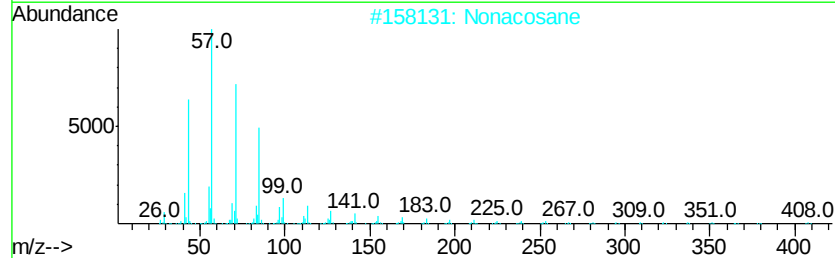
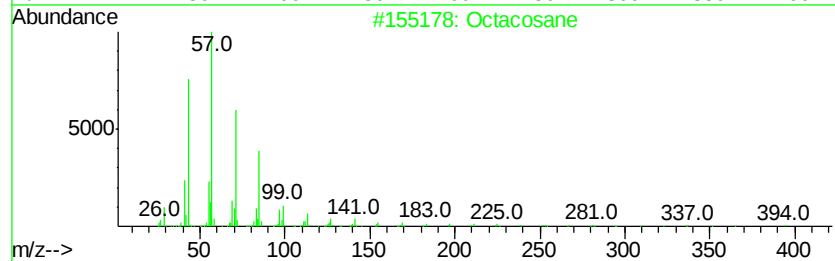
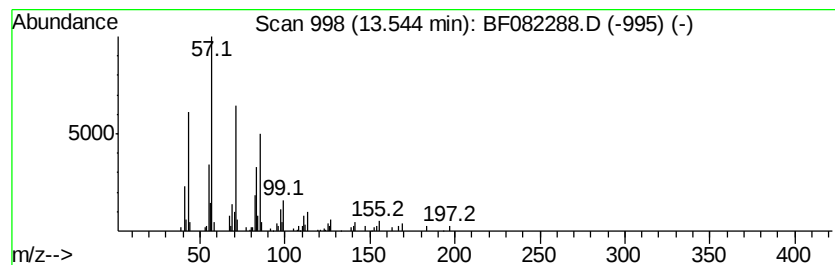
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Nonacosane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.54	3.00 ng	98617	Chrysene-d12	14.05

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octacosane	394	C28H58	000630-02-4	90
2		Nonacosane	408	C29H60	000630-03-5	87
3		Heneicosane	296	C21H44	000629-94-7	83
4		Eicosane	282	C20H42	000112-95-8	80
5		Octadecane, 1-iodo-	380	C18H37I	000629-93-6	80



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF101515\
Data File : BF082288.D
Acq On : 14 Oct 2015 19:39
Operator : UM/IZ
Sample : G4015-11
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
LF-7

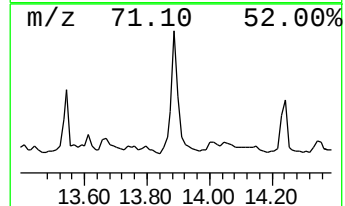
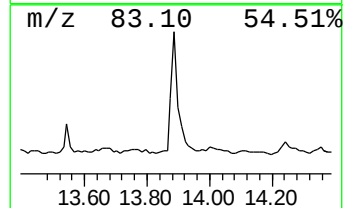
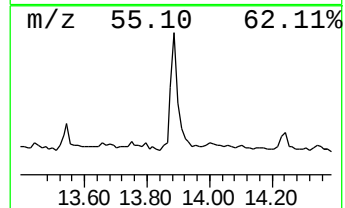
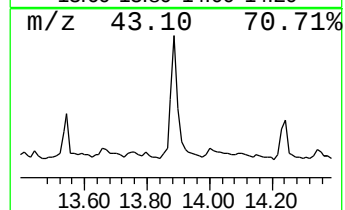
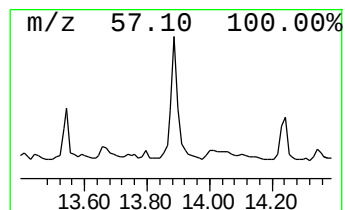
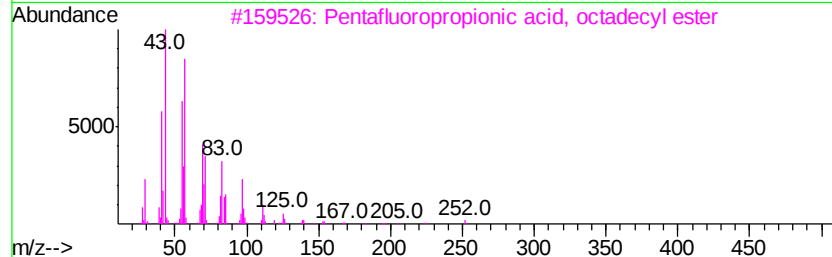
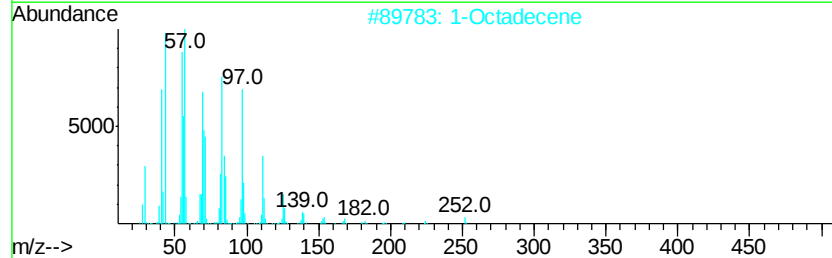
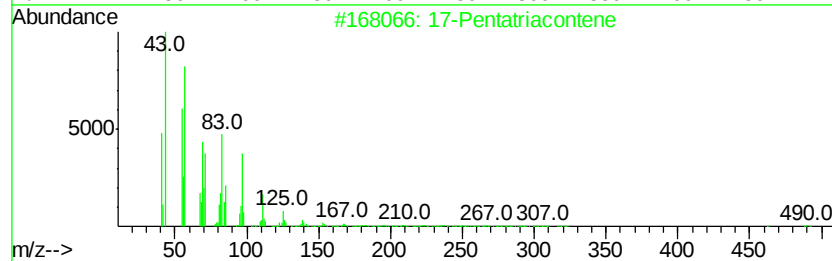
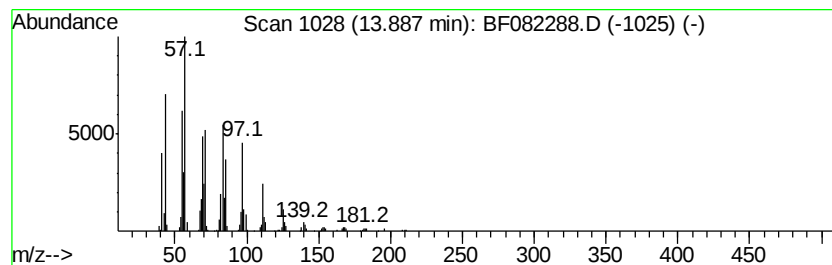
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 17-Pentatriacontene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.89	12.84 ng	422466	Chrysene-d12	14.05

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		17-Pentatriacontene	491	C35H70	006971-40-0	91
2		1-Octadecene	252	C18H36	000112-88-9	83
3		Pentafluoropropionic acid, octadecyl ester	416	C21H37F5O2	1000280-07-7	74
4		1-Heptadecanol	256	C17H36O	001454-85-9	74
5		1-Pentadecanol	228	C15H32O	000629-76-5	74



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF101515\
Data File : BF082288.D
Acq On : 14 Oct 2015 19:39
Operator : UM/IZ
Sample : G4015-11
Misc :
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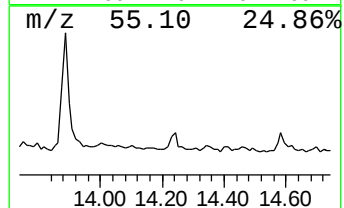
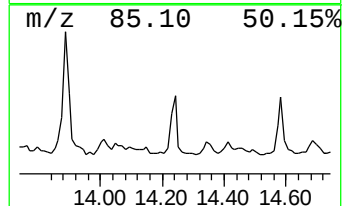
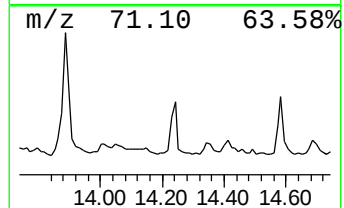
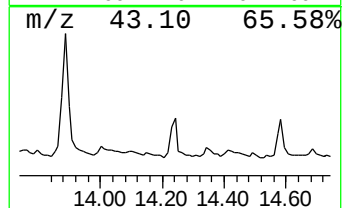
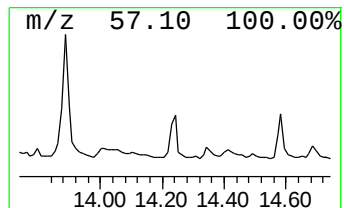
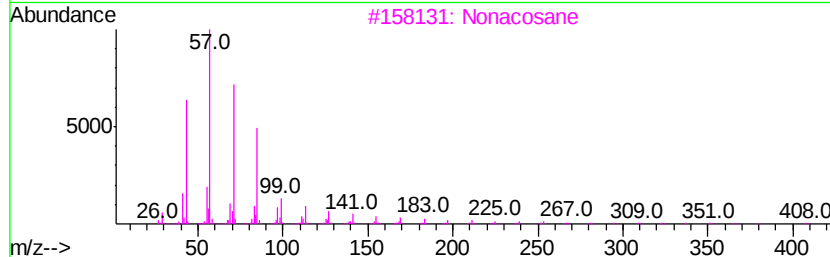
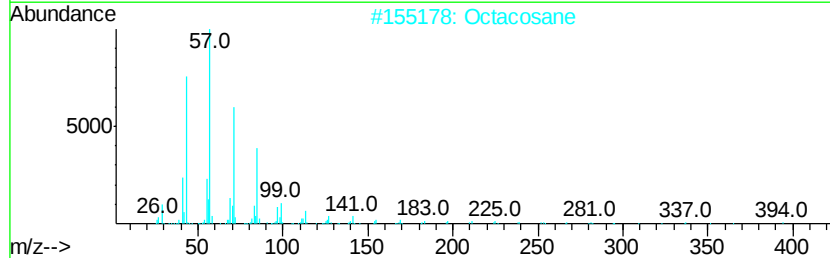
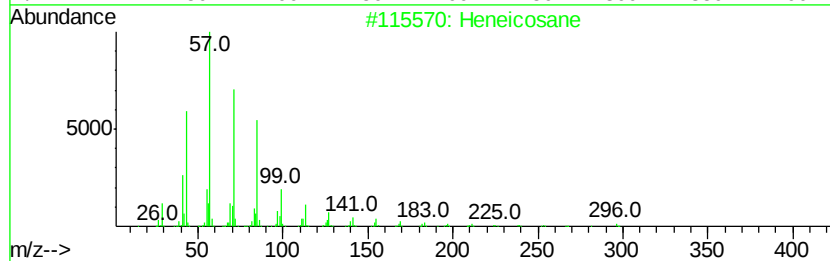
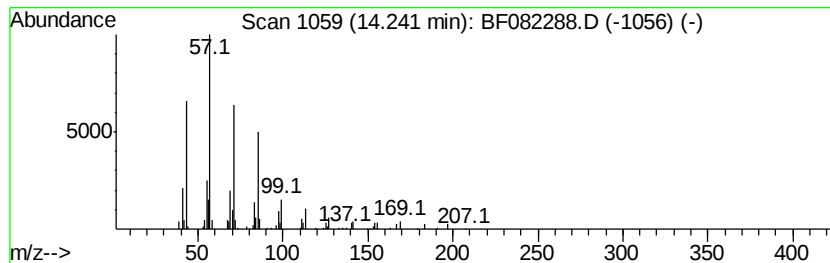
Instrument :
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF101015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Heneicosane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.24	2.46 ng	81000	Chrysene-d12	14.05
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Heneicosane	296 C21H44	000629-94-7	91
2	Octacosane	394 C28H58	000630-02-4	91
3	Nonacosane	408 C29H60	000630-03-5	90
4	Heptadecane	240 C17H36	000629-78-7	90
5	Docosane	310 C22H46	000629-97-0	90



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF101515\
Data File : BF082288.D
Acq On : 14 Oct 2015 19:39
Operator : UM/IZ
Sample : G4015-11
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
LF-7

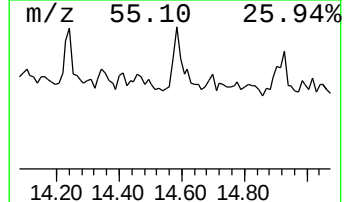
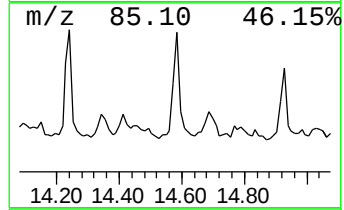
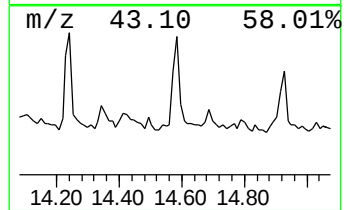
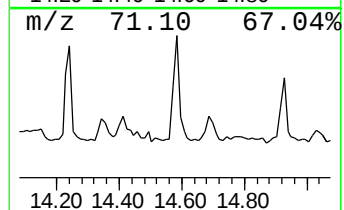
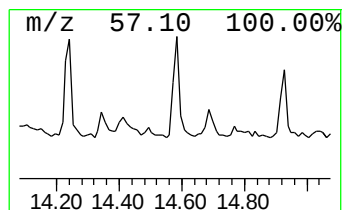
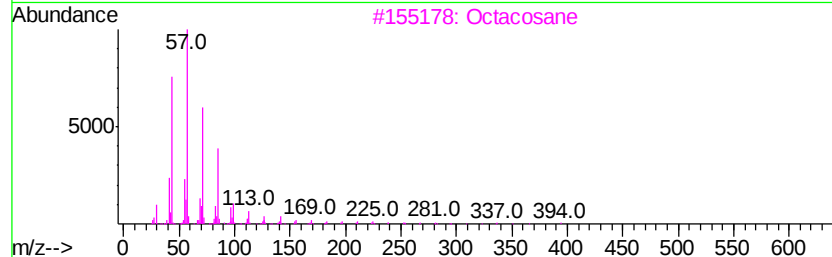
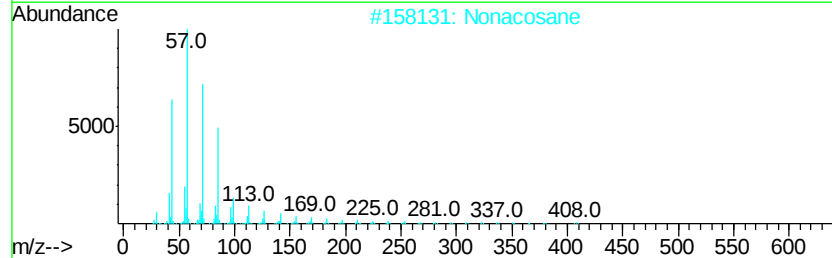
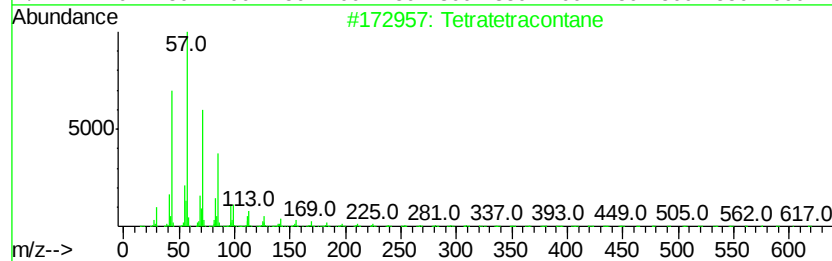
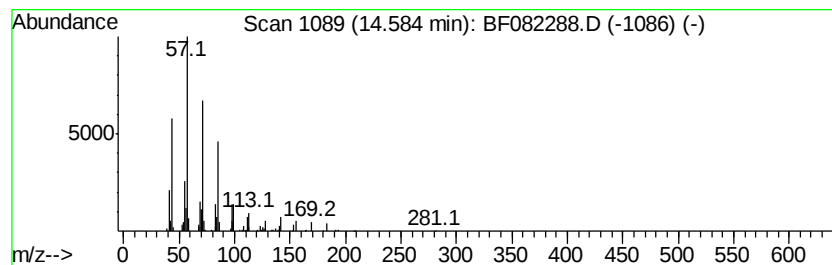
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Tetratetracontane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.58	3.39 ng	111673	Chrysene-d12	14.05

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetratetracontane	619	C44H90	007098-22-8	91
2			Nonacosane	408	C29H60	000630-03-5	91
3			Octacosane	394	C28H58	000630-02-4	87
4			Tetratriacontane	479	C34H70	014167-59-0	87
5			Docosane	310	C22H46	000629-97-0	86



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF101515\
Data File : BF082288.D
Acq On : 14 Oct 2015 19:39
Operator : UM/IZ
Sample : G4015-11
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
LF-7

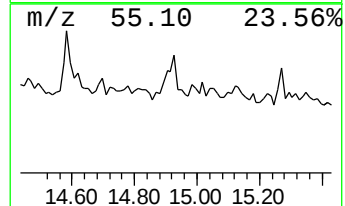
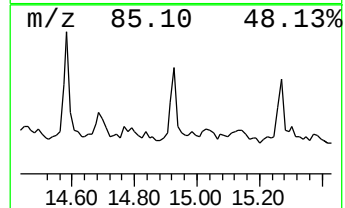
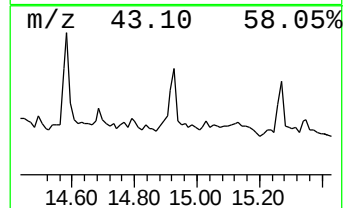
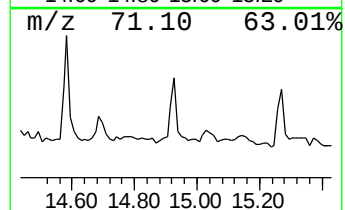
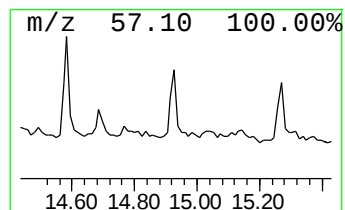
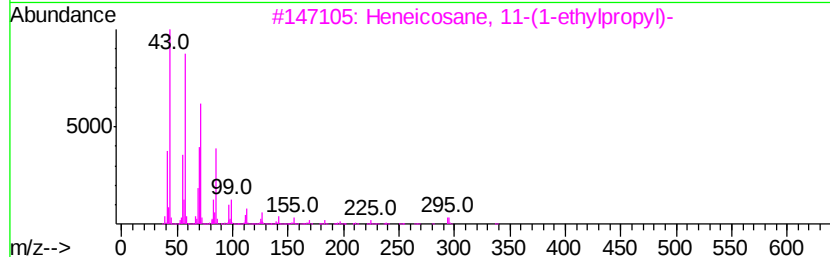
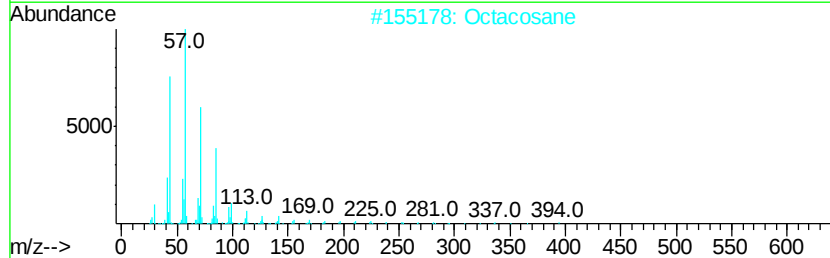
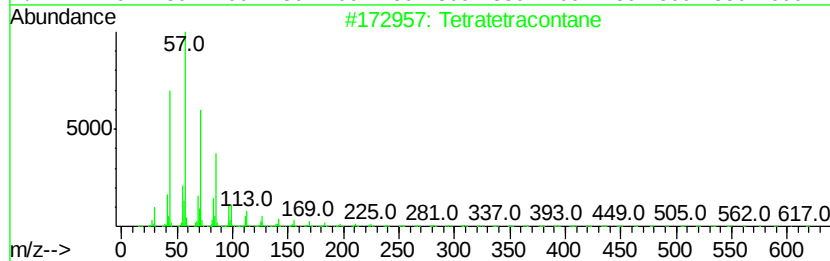
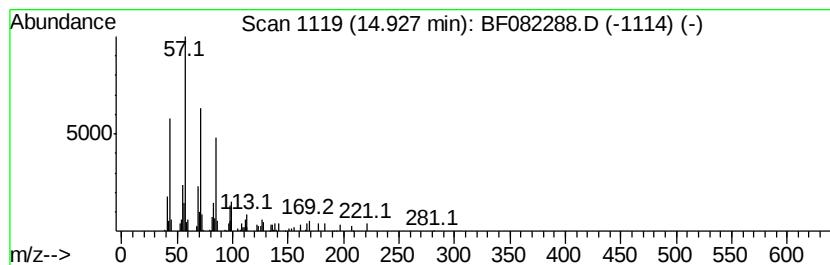
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Heneicosane, 11-(1-ethylpro... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.93	3.98 ng	97060	Perylene-d12	15.60

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetratetracontane	619	C44H90	007098-22-8	87
2			Octacosane	394	C28H58	000630-02-4	83
3			Heneicosane, 11-(1-ethylpropyl)-	366	C26H54	055282-11-6	80
4			Eicosane, 10-methyl-	296	C21H44	054833-23-7	80
5			Pentadecane, 8-heptyl-	310	C22H46	071005-15-7	80



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF101515\
Data File : BF082288.D
Acq On : 14 Oct 2015 19:39
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Sample : G4015-11
Misc :
ALS Vial : 15 Sample Multiplier: 1

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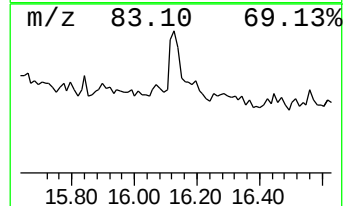
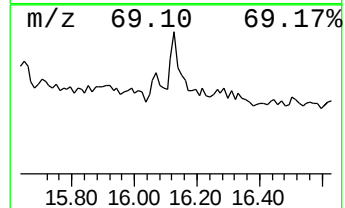
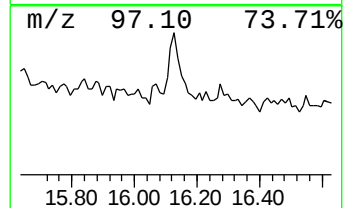
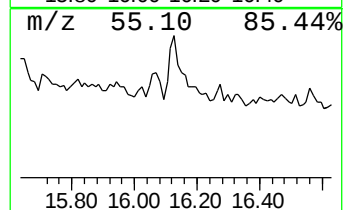
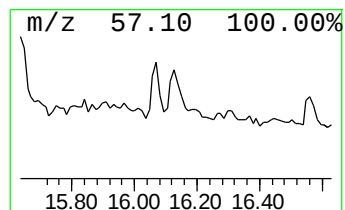
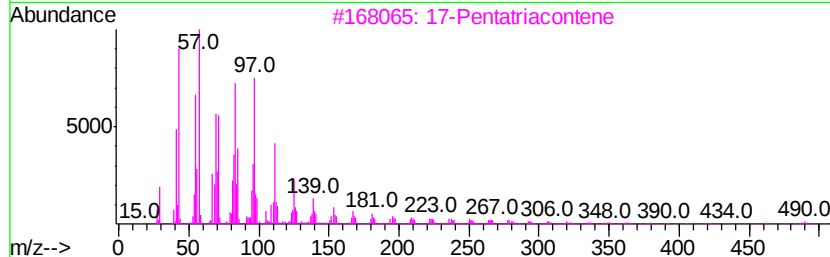
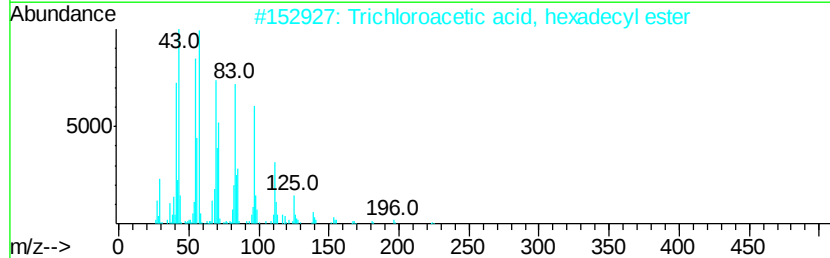
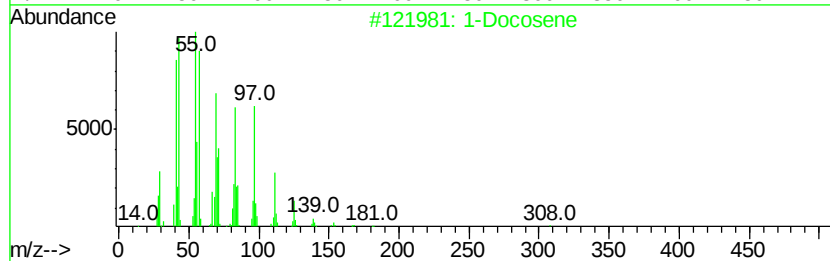
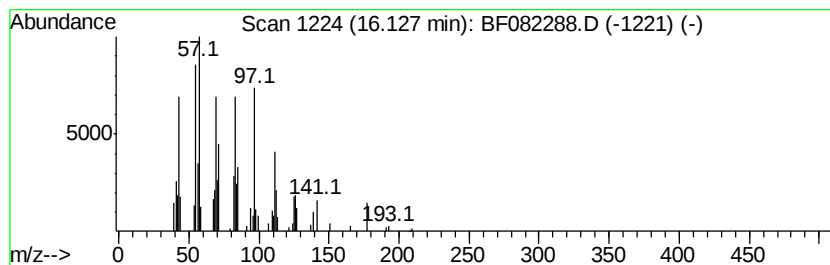
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 1-Docosene Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.13	2.02 ng	49192	Perylene-d12	15.60

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Docosene	308	C22H44	001599-67-3	80
2		Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	52
3		17-Pentatriacontene	491	C35H70	006971-40-0	49
4		Trichloroacetic acid, pentadecyl...	372	C17H31Cl3O2	074339-53-0	49
5		7-Oxabicyclo[4.1.0]heptane, 1-me...	112	C7H12O	001713-33-3	46



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF101515\
Data File : BF082288.D
Acq On : 14 Oct 2015 19:39
Operator : UM/IZ
Sample : G4015-11
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF101015.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
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2-Pentanone, 4-hy...	4.98	52.3	ng	936477	1	6.82	358037	20.0
unknown6.58	6.58	68.5	ng	1225370	1	6.82	358037	20.0
Octacosane	12.82	2.1	ng	67817	5	14.05	658090	20.0
Tridecane, 1-iodo-	13.18	2.0	ng	66710	5	14.05	658090	20.0
Nonacosane	13.54	3.0	ng	98617	5	14.05	658090	20.0
17-Pentatriacontene	13.89	12.8	ng	422466	5	14.05	658090	20.0
Heneicosane	14.24	2.5	ng	81000	5	14.05	658090	20.0
Tetratetracontane	14.58	3.4	ng	111673	5	14.05	658090	20.0
Heneicosane, 11-(...	14.93	4.0	ng	97060	6	15.60	487449	20.0
1-Docosene	16.13	2.0	ng	49192	6	15.60	487449	20.0