

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF101515\
 Data File : BF082283.D
 Acq On : 14 Oct 2015 17:01
 Operator : UM/IZ
 Sample : G4015-08
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 LF-4

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-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	4.994	247	250	252	rBV	892507	1184180	11.16%	4.060%
2	5.406	284	286	289	rBV	664441	885898	8.35%	3.037%
3	6.457	375	378	380	rBV	1362144	1301344	12.27%	4.461%
4	6.583	387	389	392	rBV	1105592	1113131	10.49%	3.816%
5	6.823	408	410	412	rBV	414896	395727	3.73%	1.357%
6	6.972	421	423	426	rBV	968886	1123765	10.59%	3.852%
7	7.394	457	460	462	rBV	817203	901838	8.50%	3.092%
8	8.114	520	523	525	rBV	446801	536824	5.06%	1.840%
9	9.189	614	617	620	rBV	1956428	1843549	17.38%	6.320%
10	9.589	649	652	654	rBV	281878	250298	2.36%	0.858%
11	9.863	674	676	679	rBV	497813	653617	6.16%	2.241%
12	10.652	743	745	750	rBV2	114806	157044	1.48%	0.538%
13	10.766	753	755	760	rVB	88760	131803	1.24%	0.452%
14	11.361	804	807	809	rBV	628137	697145	6.57%	2.390%
15	11.601	824	828	830	rBV3	59198	142067	1.34%	0.487%
16	11.646	830	832	835	rVV	129403	126183	1.19%	0.433%
17	11.829	844	848	851	rBV4	54146	192748	1.82%	0.661%
18	11.909	851	855	857	rBV3	63493	161127	1.52%	0.552%
19	12.058	867	868	870	rBV	171361	172931	1.63%	0.593%
20	12.344	891	893	895	rBV	146931	202894	1.91%	0.696%
21	12.412	897	899	901	rBV2	104058	164716	1.55%	0.565%
22	12.458	901	903	904	rBV	176932	211054	1.99%	0.724%
23	12.641	918	919	920	rVB	156346	107253	1.01%	0.368%
24	12.664	920	921	923	rBV	126366	222414	2.10%	0.762%
25	12.835	935	936	938	rBV	281216	297531	2.80%	1.020%
26	12.961	944	947	950	rBV	1594924	2095750	19.75%	7.185%
27	13.201	966	968	970	rBV2	259882	309478	2.92%	1.061%
28	13.566	998	1000	1002	rBV	335911	376253	3.55%	1.290%
29	14.092	1041	1046	1049	rBV	4968818	10609514	100.00%	36.371%
30	14.618	1090	1092	1095	rBV	679048	859774	8.10%	2.947%
31	15.658	1181	1183	1186	rVB	302277	388139	3.66%	1.331%
32	16.344	1241	1243	1251	rVB3	257557	713694	6.73%	2.447%
33	16.767	1277	1280	1288	rVB2	238119	640451	6.04%	2.196%

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Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

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

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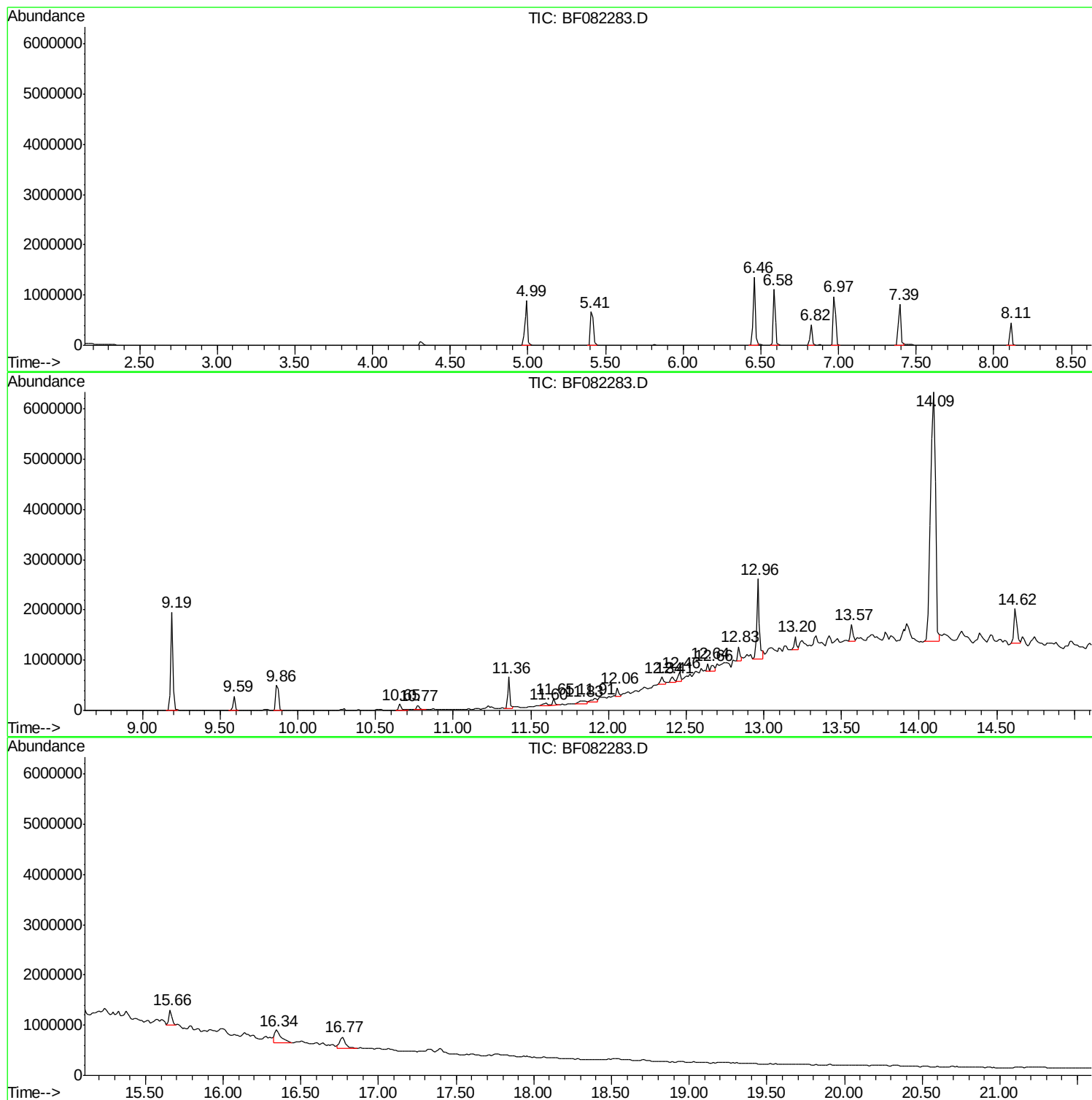
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF101015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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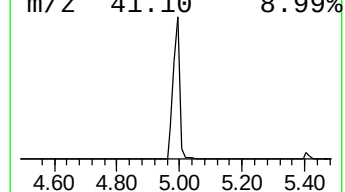
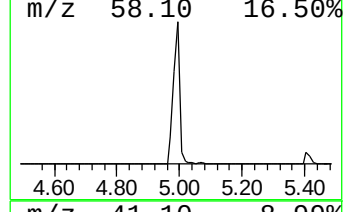
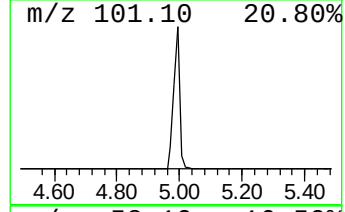
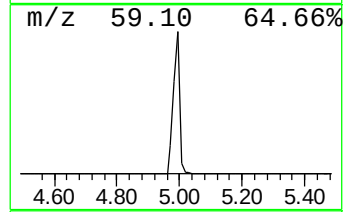
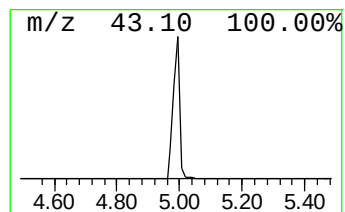
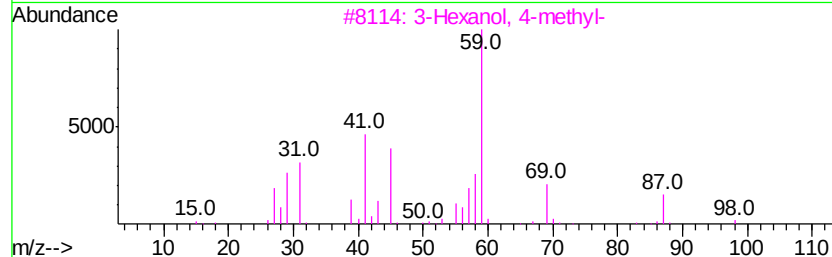
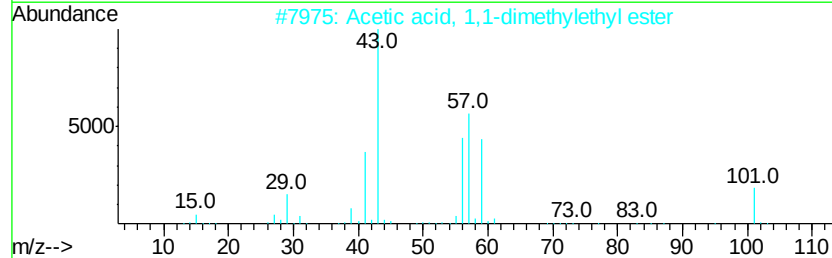
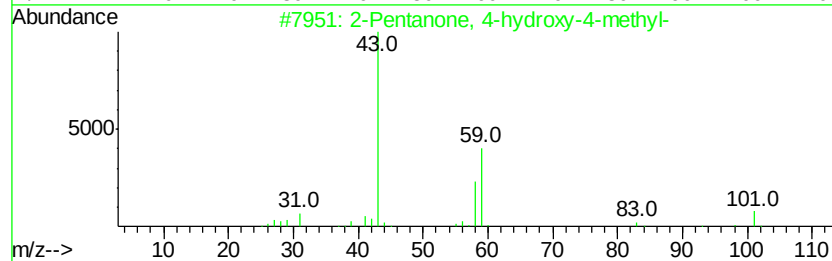
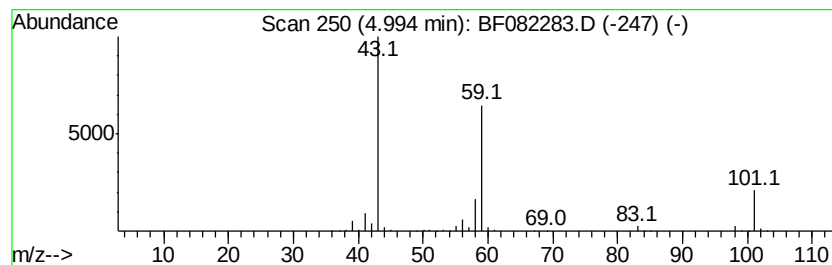
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.99	59.85 ng	1184180	1,4-Dichlorobenzene-d4	6.82

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2		Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
3		3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	28
4		Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	28
5		Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	17



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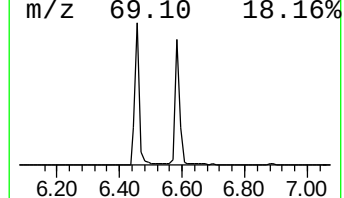
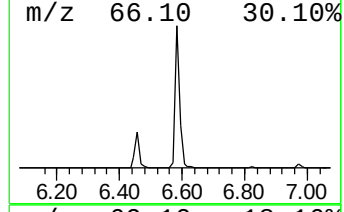
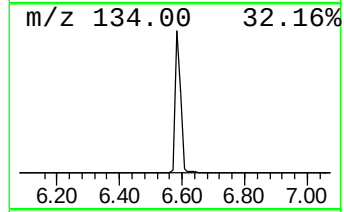
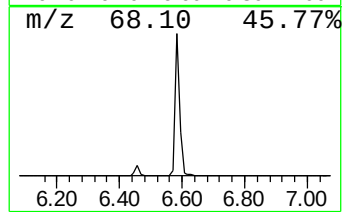
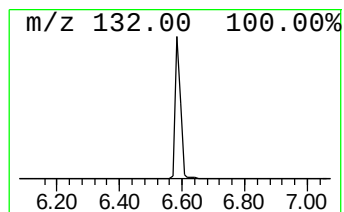
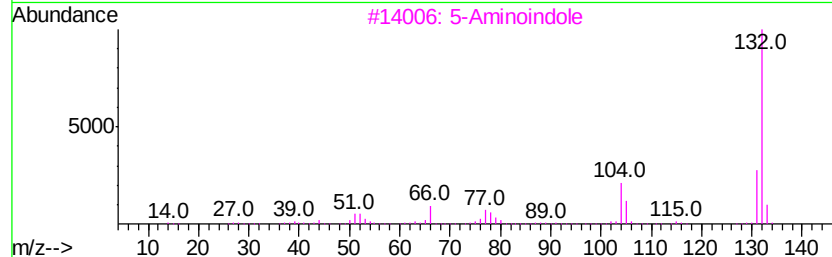
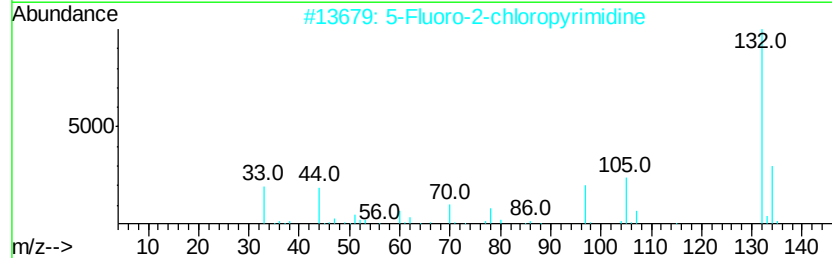
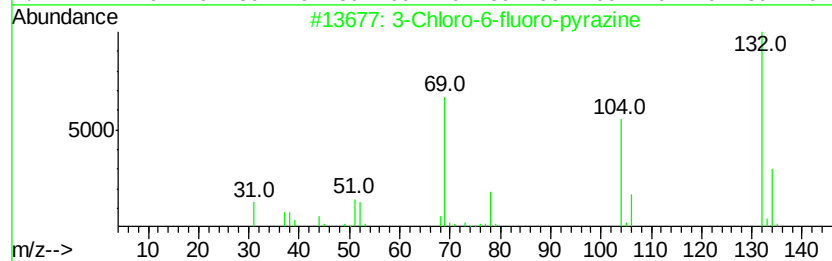
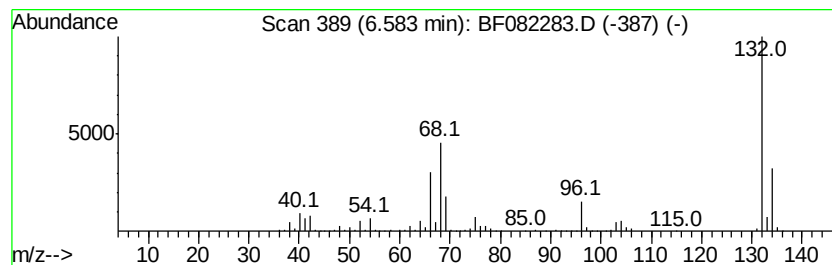
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 2 unknown6.58 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.58	56.26 ng	1113130	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	27
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
3		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
5		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9



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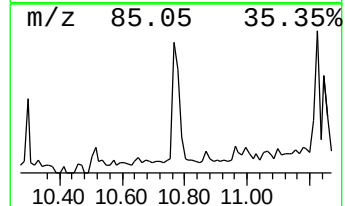
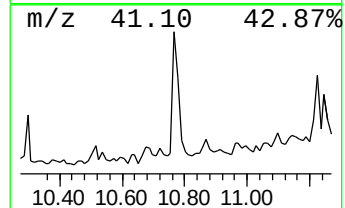
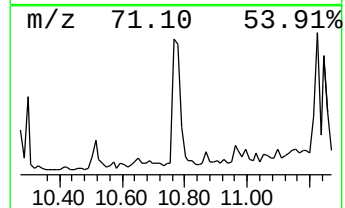
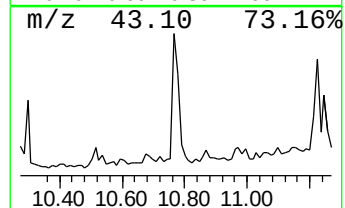
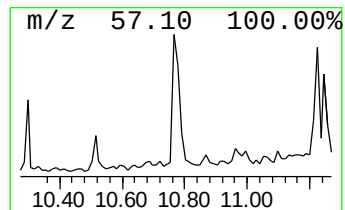
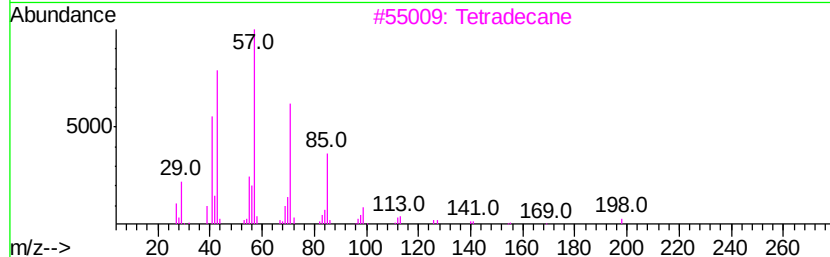
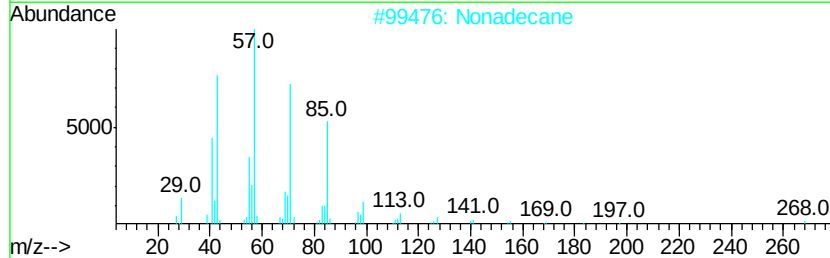
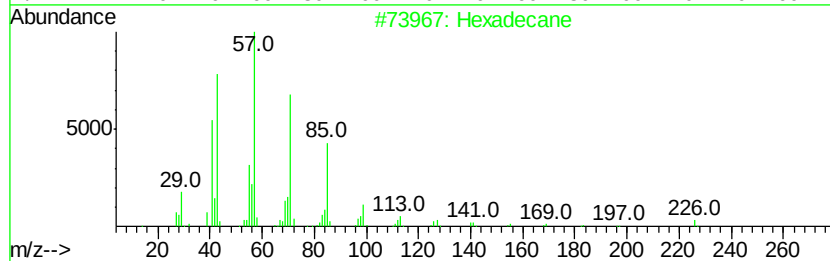
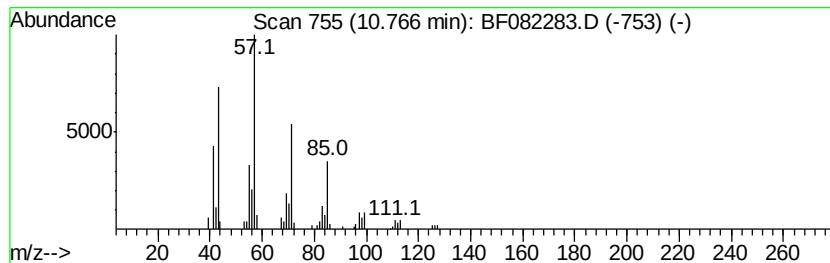
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Hexadecane Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.77	3.78 ng	131803	Phenanthrene-d10	11.36	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane	226	C16H34	000544-76-3	86
2	Nonadecane	268	C19H40	000629-92-5	86
3	Tetradecane	198	C14H30	000629-59-4	86
4	Dotriacontane	451	C32H66	000544-85-4	86
5	1-Iodo-2-methylundecane	296	C12H25I	073105-67-6	78



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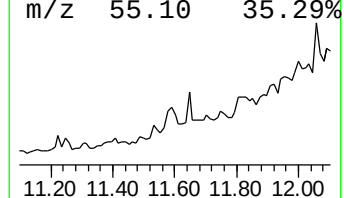
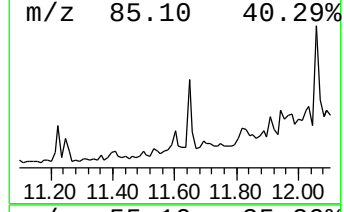
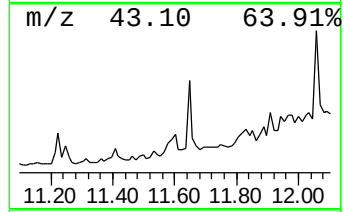
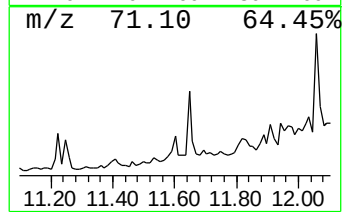
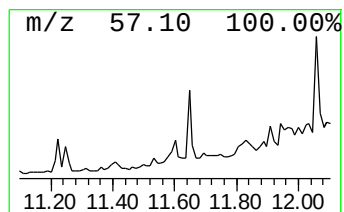
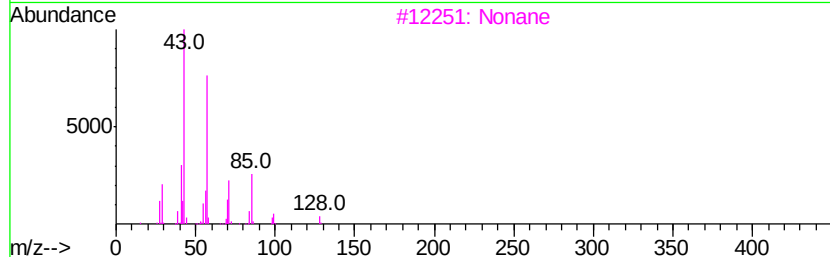
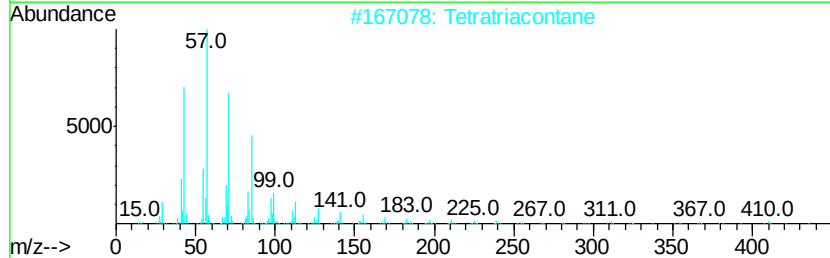
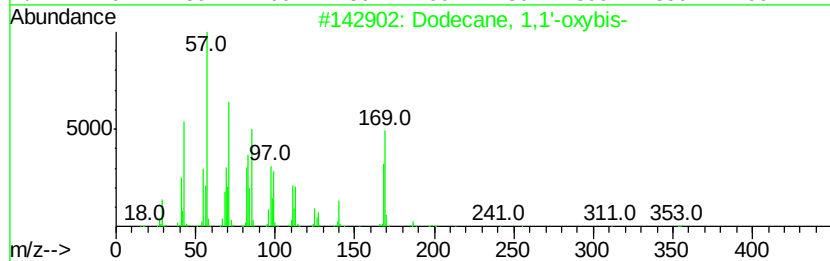
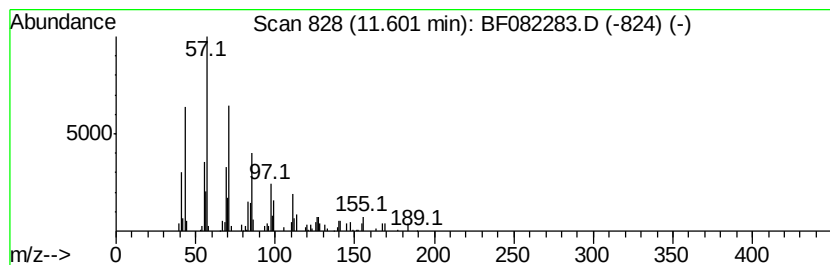
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF101015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Dodecane, 1,1'-oxybis- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.60	4.08 ng	142067	Phenanthrene-d10	11.36

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Dodecane, 1,1'-oxybis-	354	C24H50O	004542-57-8	83
2		Tetratriacontane	479	C34H70	014167-59-0	64
3		Nonane	128	C9H20	000111-84-2	60
4		Tetratetracontane	619	C44H90	007098-22-8	59
5		Tetradecane, 2,6,10-trimethyl-	240	C17H36	014905-56-7	59



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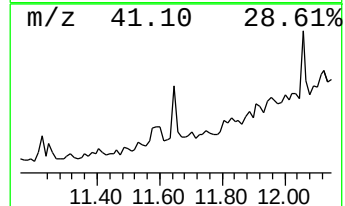
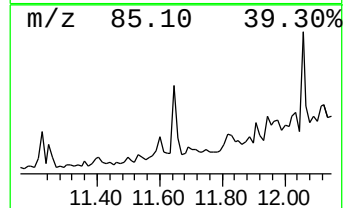
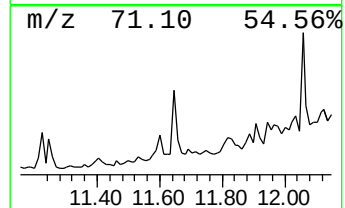
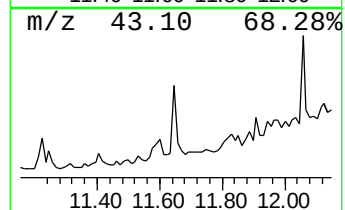
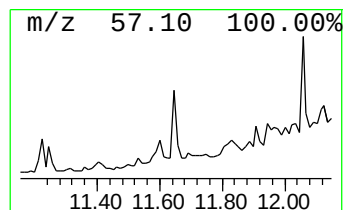
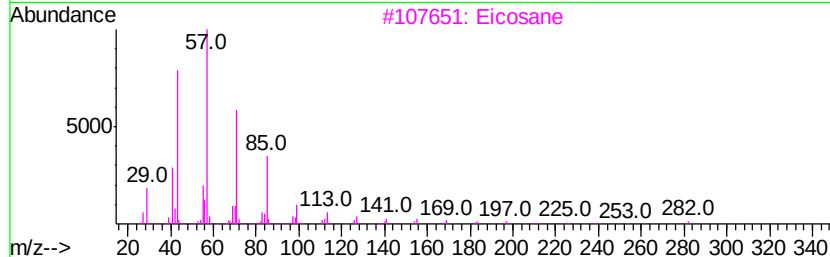
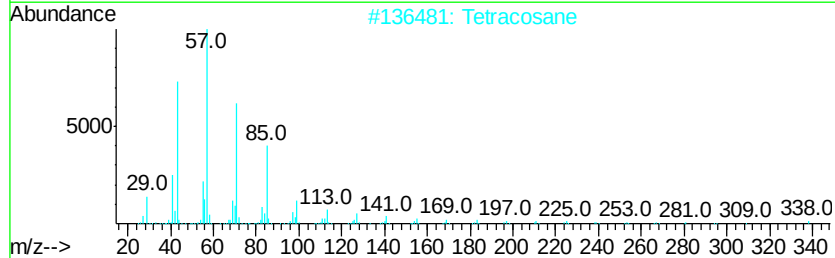
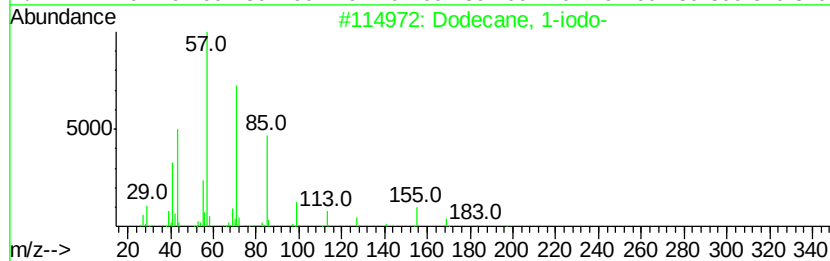
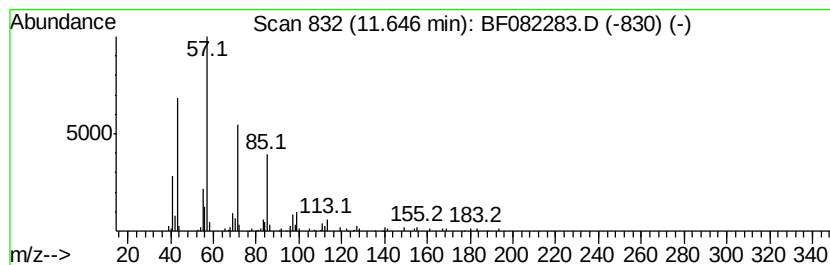
Instrument :
BNA_F
ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Dodecane, 1-iodo- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.65	3.62 ng	126183	Phenanthrene-d10	11.36	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dodecane, 1-iodo-	296	C12H25I	004292-19-7	86
2	Tetracosane	338	C24H50	000646-31-1	86
3	Eicosane	282	C20H42	000112-95-8	80
4	Eicosane, 10-methyl-	296	C21H44	054833-23-7	80
5	Octadecane	254	C18H38	000593-45-3	72



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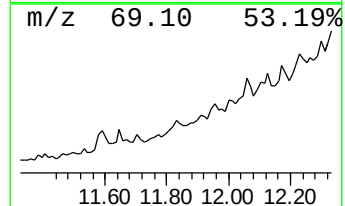
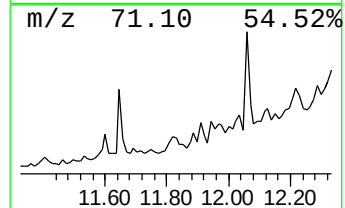
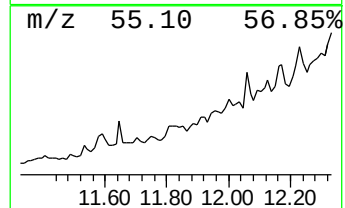
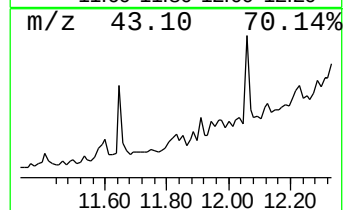
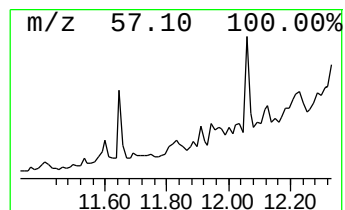
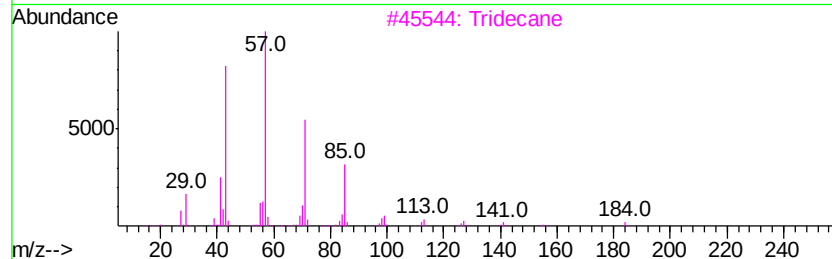
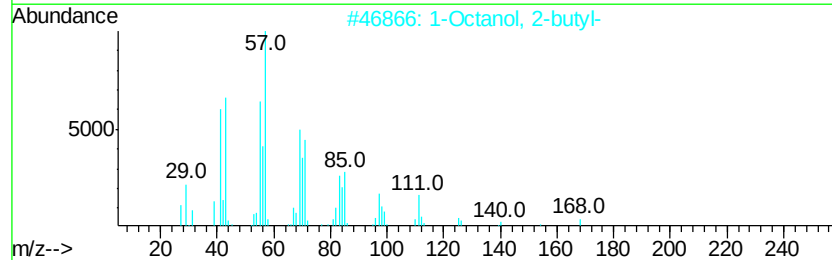
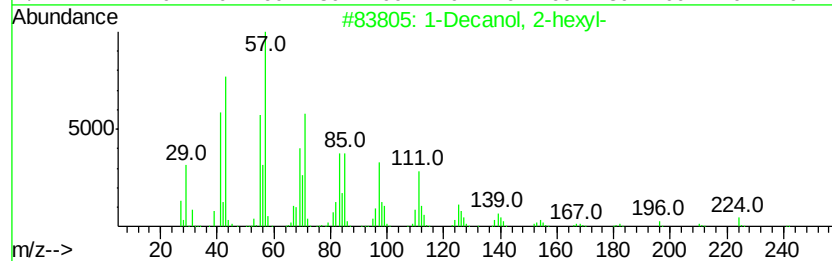
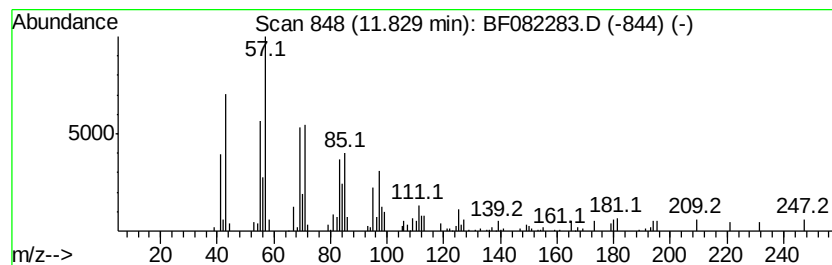
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF101015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 1-Decanol, 2-hexyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.83	5.53 ng	192748	Phenanthrene-d10	11.36	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Decanol, 2-hexyl-	242	C16H34O	002425-77-6	62
2	1-Octanol, 2-butyl-	186	C12H26O	003913-02-8	58
3	Tridecane	184	C13H28	000629-50-5	43
4	Dichloroacetic acid, 2-pentadecyl...	338	C17H32Cl2O2	1000280-64-7	43
5	Heptadecane	240	C17H36	000629-78-7	43



Data Path : Z:\HPCHEM1\BNA F\DATA\BF101515\
Data File : BF082283.D
Acq On : 14 Oct 2015 17:01
Operator : UM/IZ
Sample : G4015-08
Misc :
ALS Vial : 10 Sample Multiplier: 1

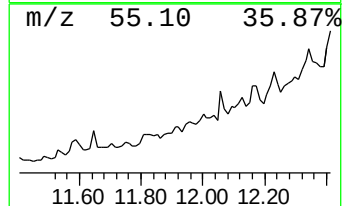
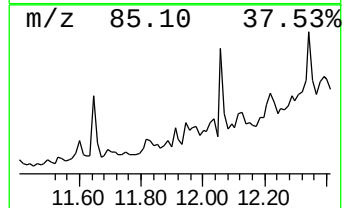
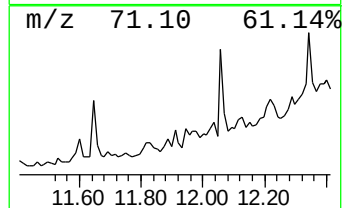
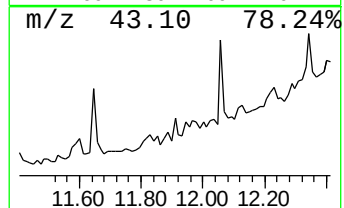
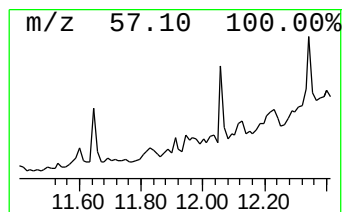
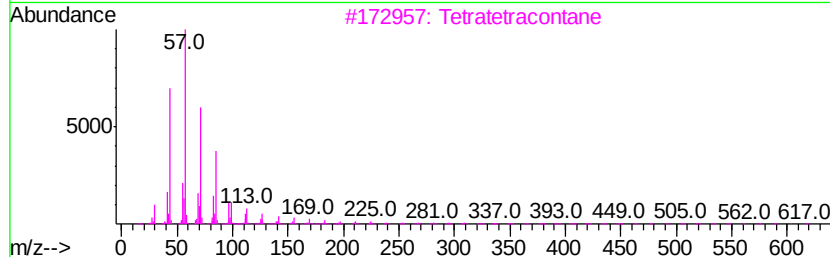
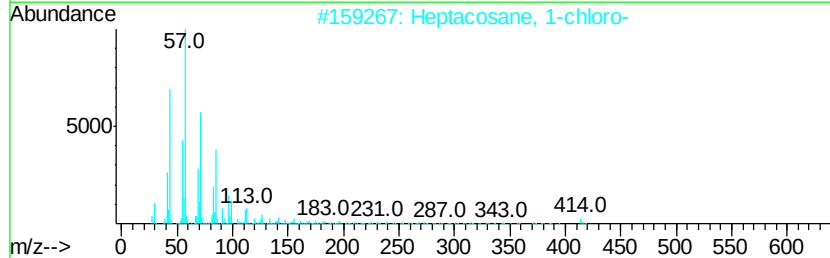
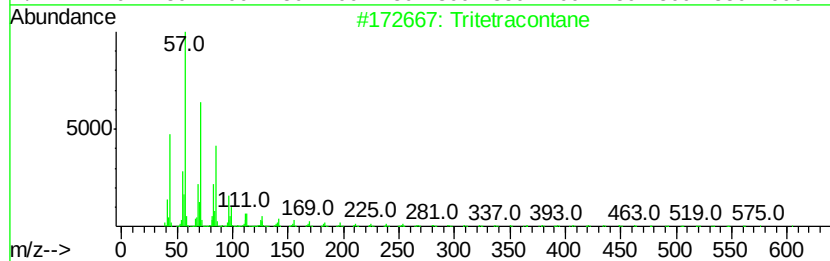
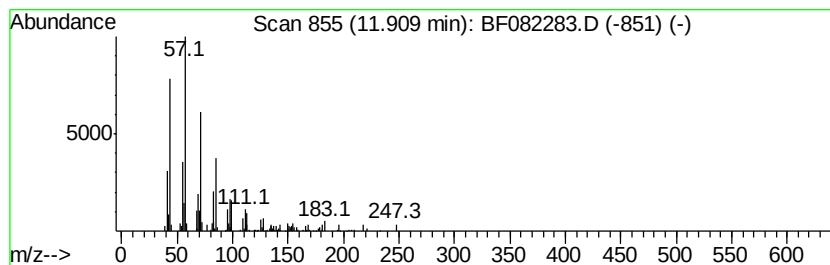
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF101015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Tritetracontane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD		R.T.
11.91	4.62 ng	161127	Phenanthrene-d10		11.36
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tritetracontane	605	C43H88	007098-21-7	80
2	Heptacosane, 1-chloro-	414	C27H55Cl	062016-79-9	80
3	Tetratetracontane	619	C44H90	007098-22-8	80
4	Heptacosane	380	C27H56	000593-49-7	64
5	Octacosane	394	C28H58	000630-02-4	59



Data Path : Z:\HPCHEM1\BNA F\DATA\BF101515\
Data File : BF082283.D
Acq On : 14 Oct 2015 17:01
Operator : UM/IZ
Sample : G4015-08
Misc :
ALS Vial : 10 Sample Multiplier: 1

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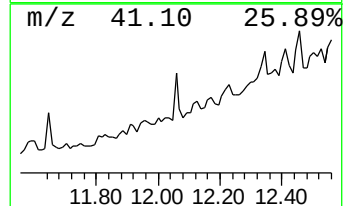
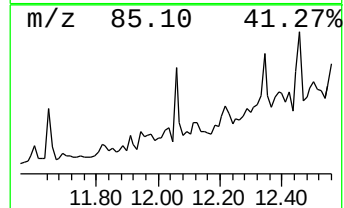
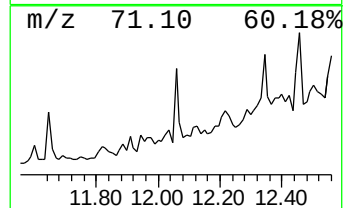
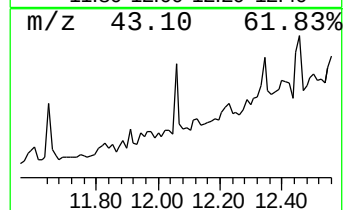
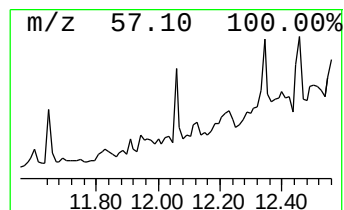
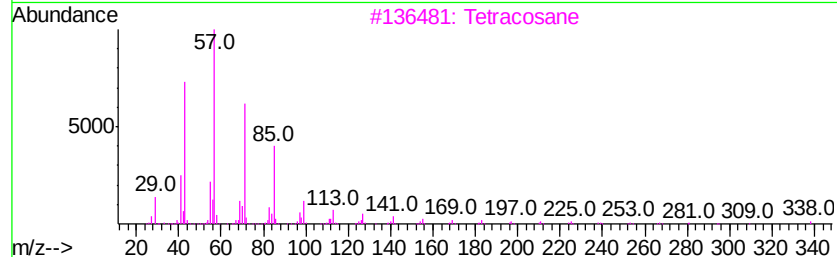
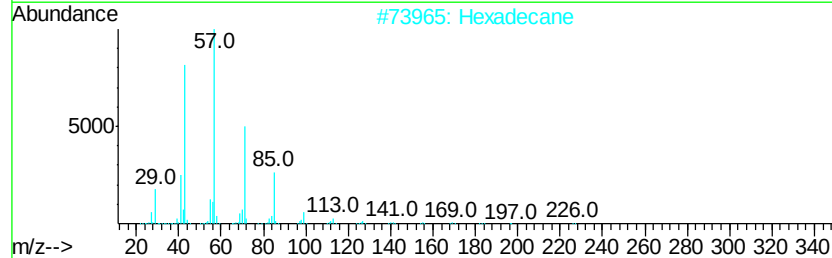
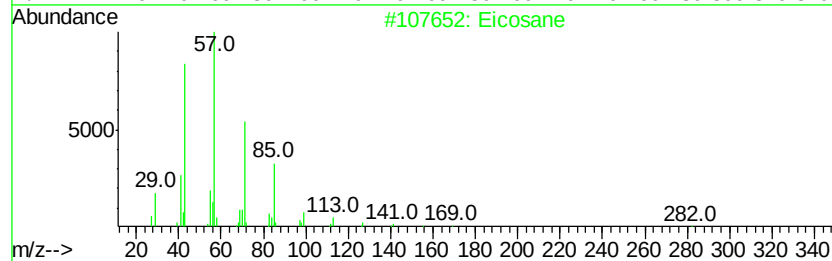
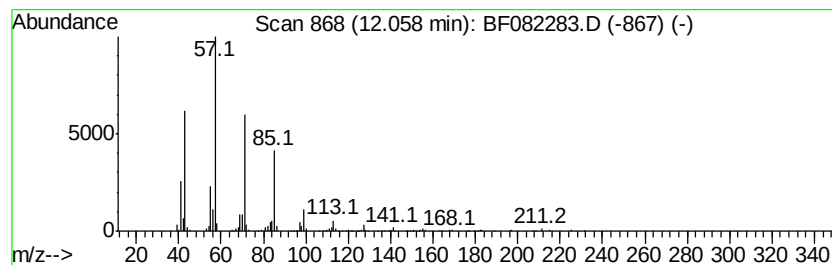
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Eicosane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.06	4.96 ng	172931	Phenanthrene-d10	11.36

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane	282	C20H42	000112-95-8	91
2	Hexadecane	226	C16H34	000544-76-3	90
3	Tetracosane	338	C24H50	000646-31-1	90
4	Tetradecane, 1-iodo-	324	C14H29I	019218-94-1	86
5	Dodecane, 1-iodo-	296	C12H25I	004292-19-7	86



Data Path : Z:\HPCHEM1\BNA F\DATA\BF101515\
Data File : BF082283.D
Acq On : 14 Oct 2015 17:01
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Sample : G4015-08
Misc :
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Instrument :
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ClientSampleId :
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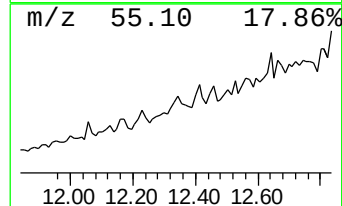
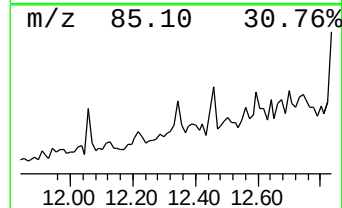
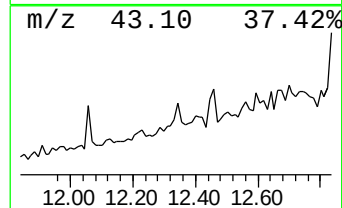
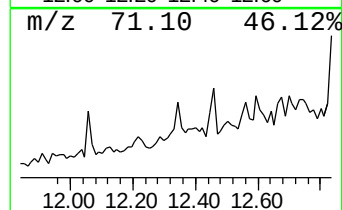
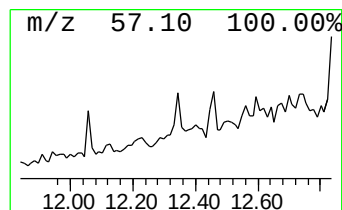
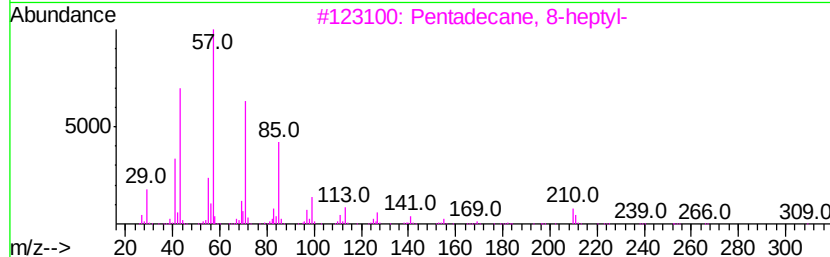
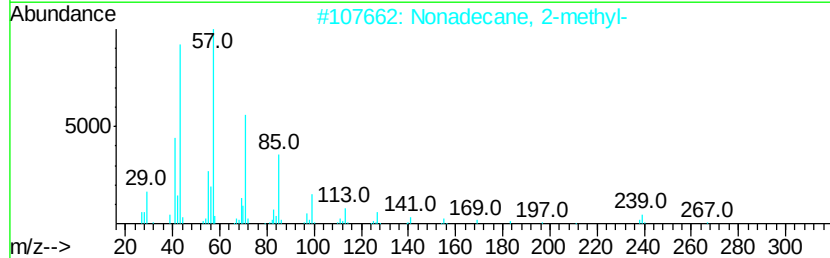
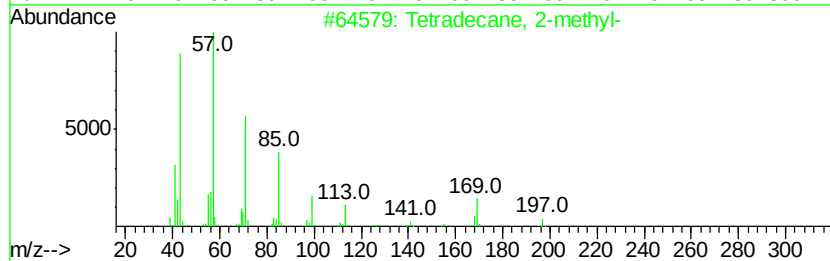
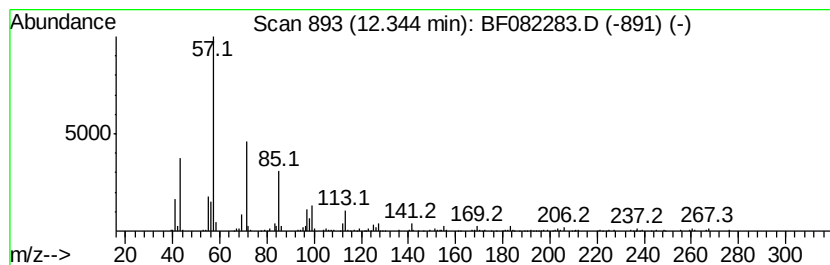
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Tetradecane, 2-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.34	5.82 ng	202894	Phenanthrene-d10	11.36

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetradecane, 2-methyl-	212	C15H32	001560-95-8	87
2			Nonadecane, 2-methyl-	282	C20H42	001560-86-7	86
3			Pentadecane, 8-heptyl-	310	C22H46	071005-15-7	86
4			Tetradecane, 3-methyl-	212	C15H32	018435-22-8	83
5			Heptadecane	240	C17H36	000629-78-7	80



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF101515\
Data File : BF082283.D
Acq On : 14 Oct 2015 17:01
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Sample : G4015-08
Misc :
ALS Vial : 10 Sample Multiplier: 1

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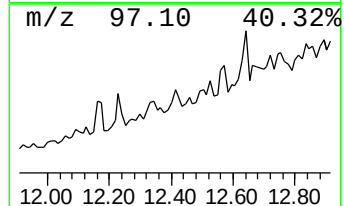
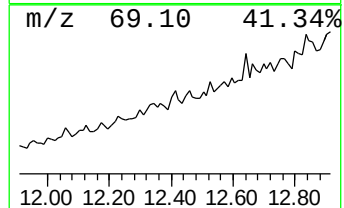
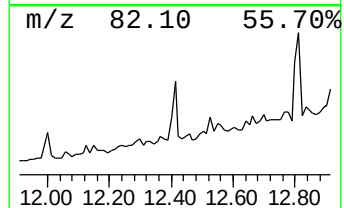
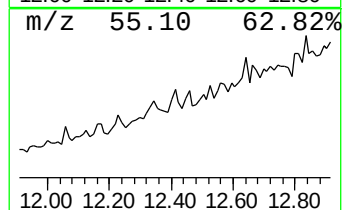
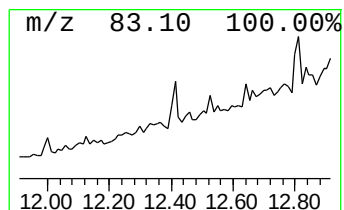
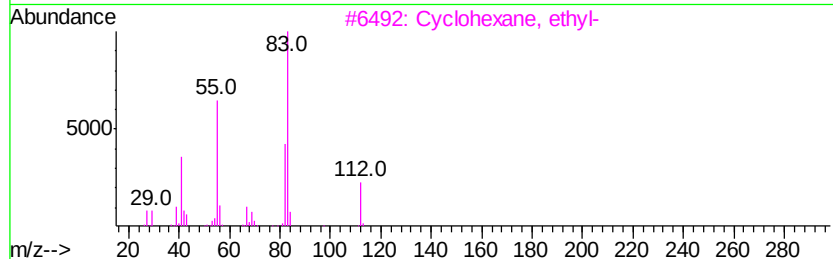
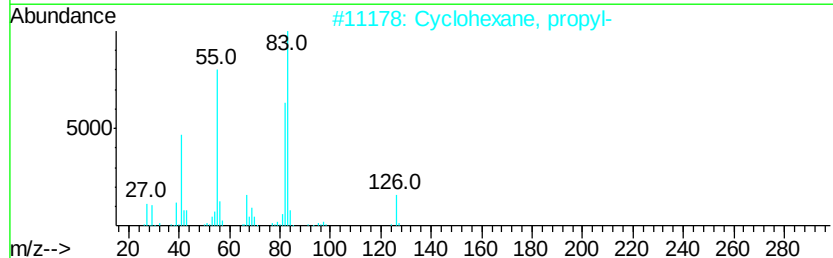
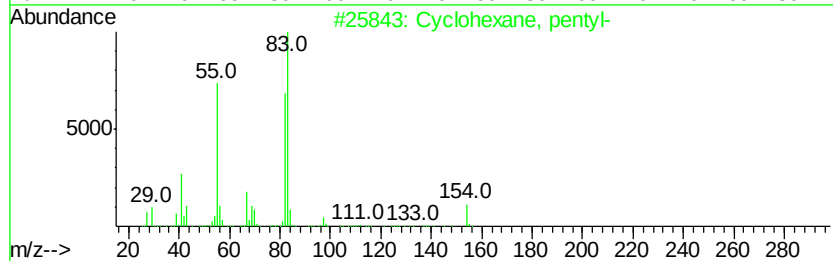
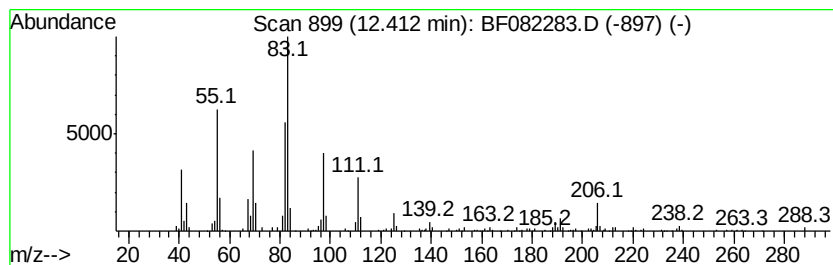
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 unknown12.41 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.41	4.73 ng	164716	Phenanthrene-d10	11.36

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, pentyl-	154	C11H22	004292-92-6	47
2			Cyclohexane, propyl-	126	C9H18	001678-92-8	47
3			Cyclohexane, ethyl-	112	C8H16	001678-91-7	47
4			Cyclohexanecarbonyl chloride	146	C7H11ClO	002719-27-9	38
5			Cyclohexane carboxylic acid hydr...	142	C7H14N2O	038941-47-8	38



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF101515\
Data File : BF082283.D
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Operator : UM/IZ
Sample : G4015-08
Misc :
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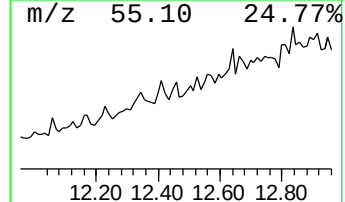
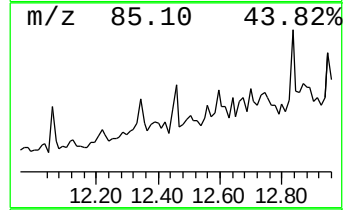
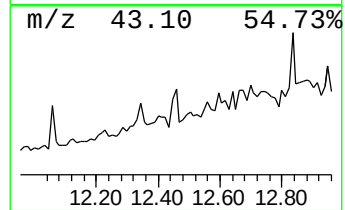
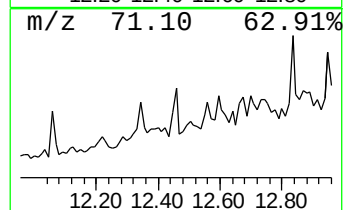
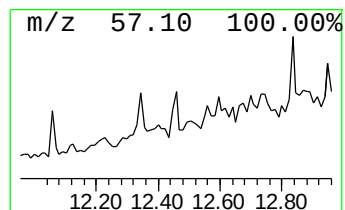
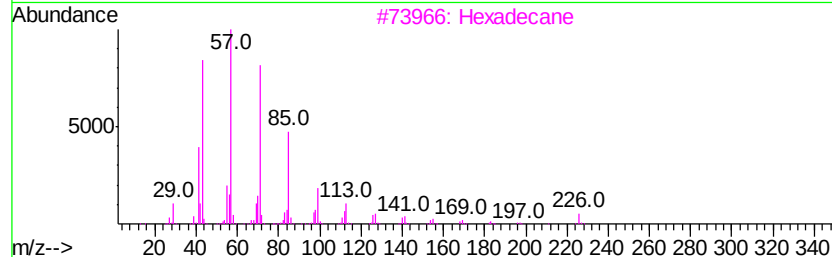
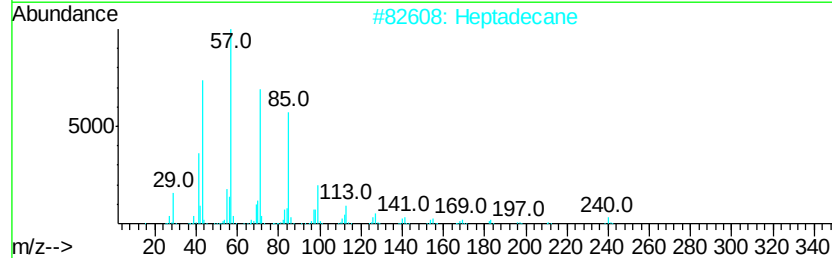
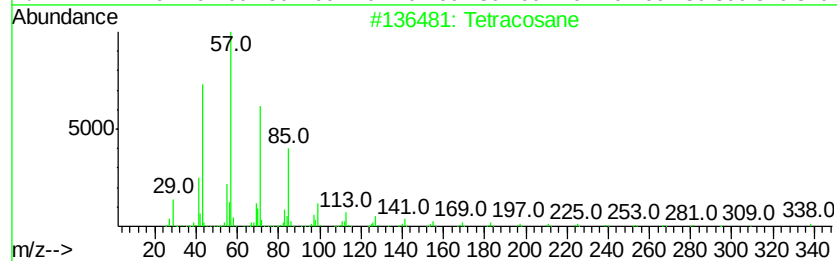
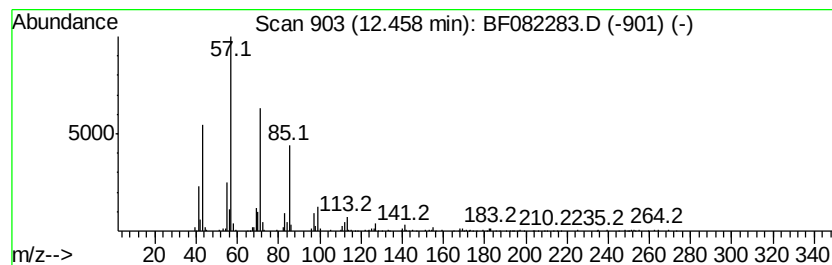
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Tetracosane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD		R.T.
12.46	6.05 ng	211054	Phenanthrene-d10		11.36
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetracosane	338	C24H50	000646-31-1	87
2	Heptadecane	240	C17H36	000629-78-7	86
3	Hexadecane	226	C16H34	000544-76-3	86
4	Eicosane	282	C20H42	000112-95-8	86
5	Octadecane, 2-methyl-	268	C19H40	001560-88-9	86



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF101515\
Data File : BF082283.D
Acq On : 14 Oct 2015 17:01
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Instrument :
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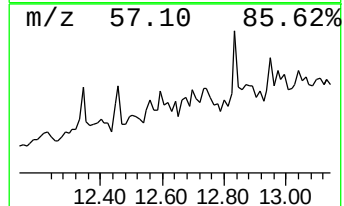
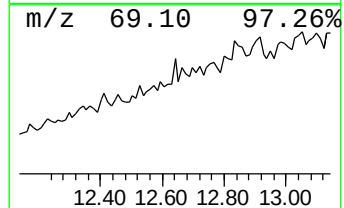
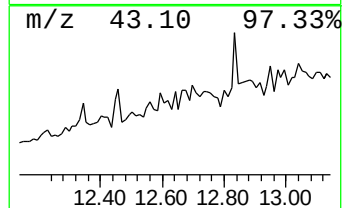
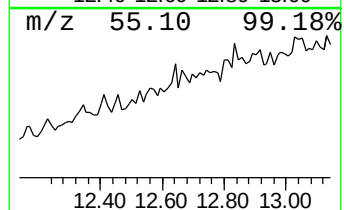
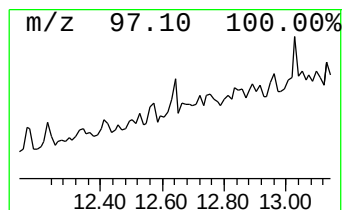
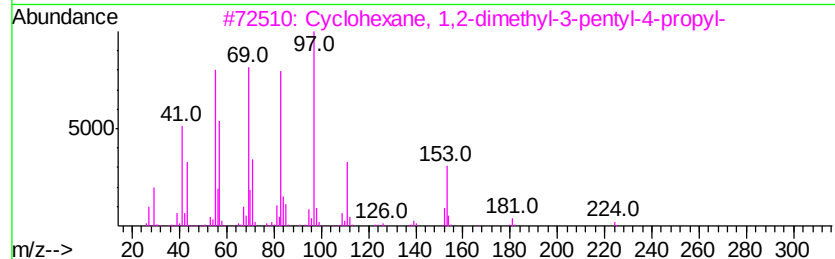
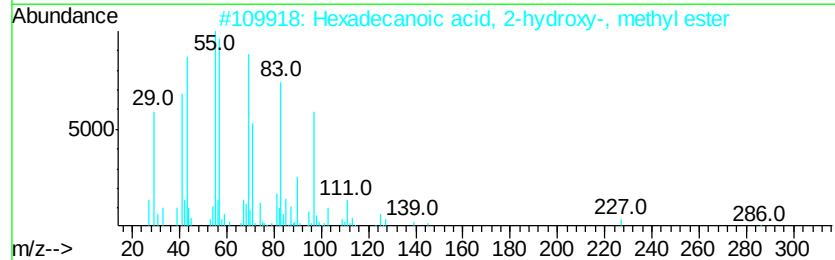
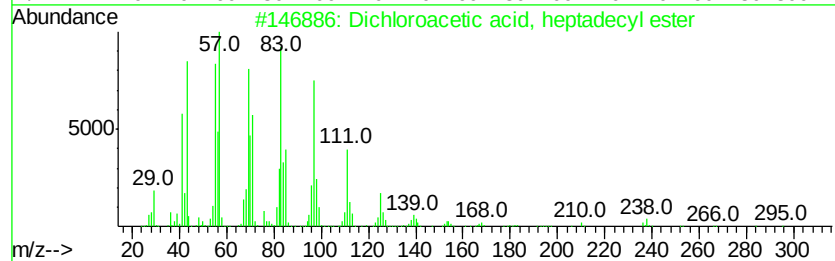
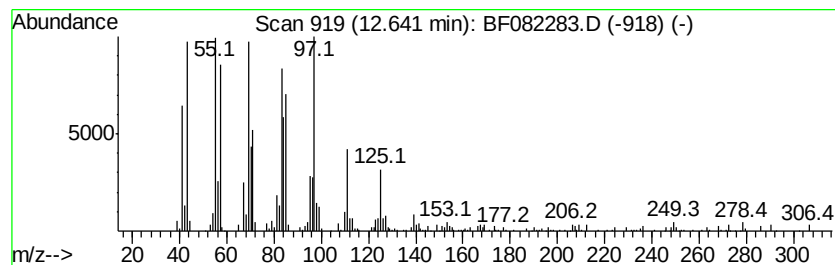
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Dichloroacetic acid, heptad... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.64	3.08 ng	107253	Phenanthrene-d10	11.36

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dichloroacetic acid, heptadecyl ...	366	C19H36Cl2O2	1000282-98-2	80
2			Hexadecanoic acid, 2-hydroxy-, m...	286	C17H34O3	016742-51-1	70
3			Cyclohexane, 1,2-dimethyl-3-pent...	224	C16H32	062376-17-4	70
4			Bromoacetic acid, octadecyl ester	390	C20H39BrO2	018992-03-5	59
5			2-Cyclohexen-1-ol, 1-methyl-	112	C7H12O	023758-27-2	41



Data Path : Z:\HPCHEM1\BNA F\DATA\BF101515\
Data File : BF082283.D
Acq On : 14 Oct 2015 17:01
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Sample : G4015-08
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
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ClientSampleId :
LF-4

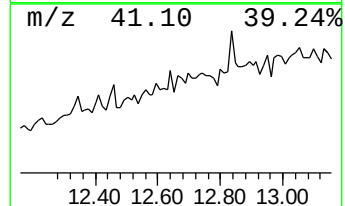
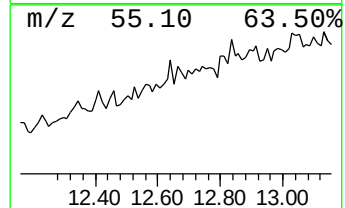
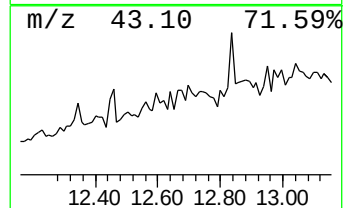
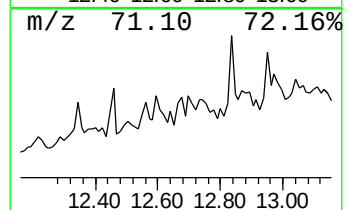
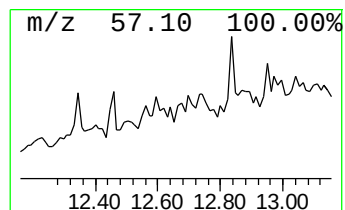
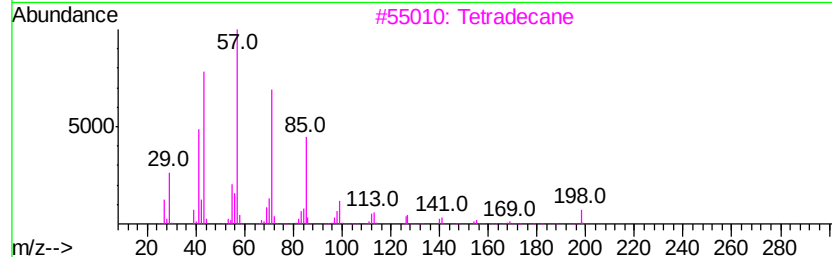
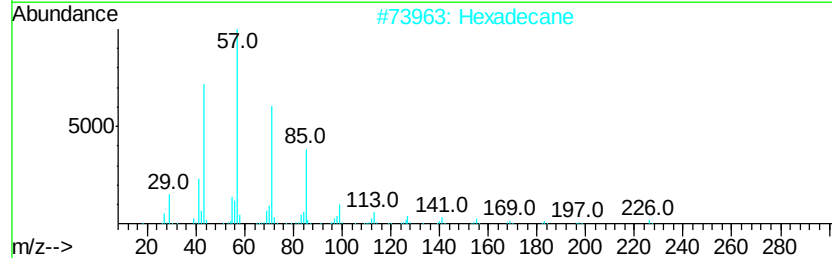
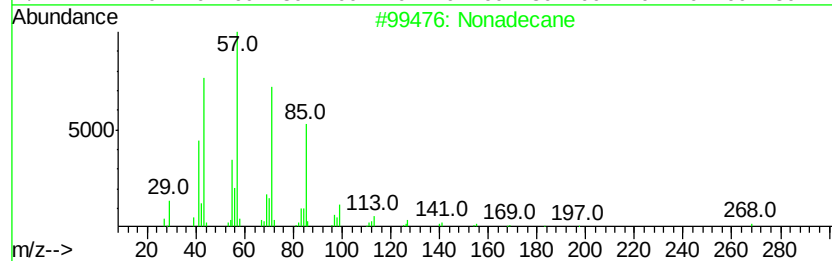
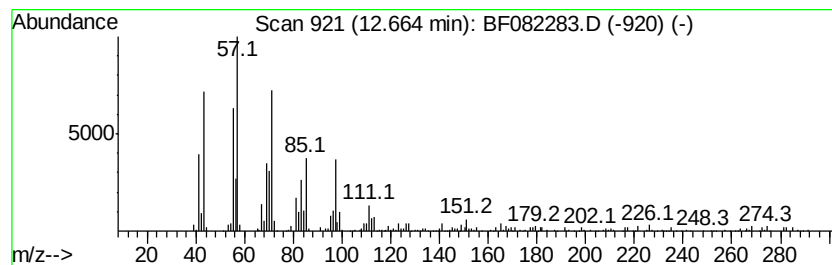
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Nonadecane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.66	6.38 ng	222414	Phenanthrene-d10	11.36

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonadecane	268	C19H40	000629-92-5	91
2			Hexadecane	226	C16H34	000544-76-3	90
3			Tetradecane	198	C14H30	000629-59-4	78
4			Eicosane	282	C20H42	000112-95-8	78
5			Tritetracontane	605	C43H88	007098-21-7	72



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF101515\
Data File : BF082283.D
Acq On : 14 Oct 2015 17:01
Operator : UM/IZ
Sample : G4015-08
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
LF-4

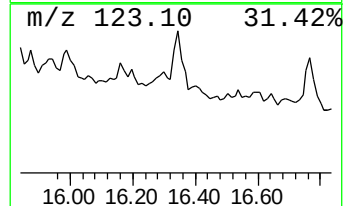
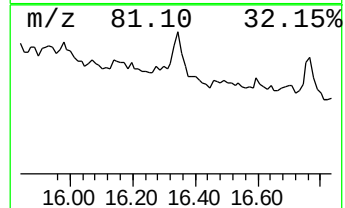
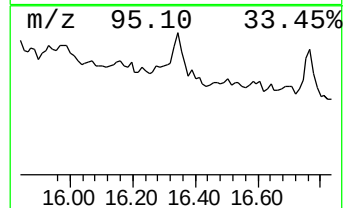
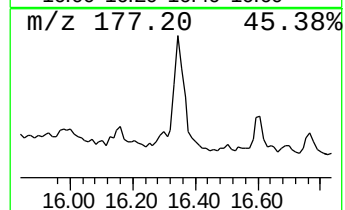
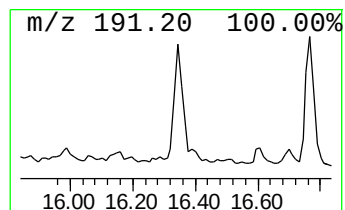
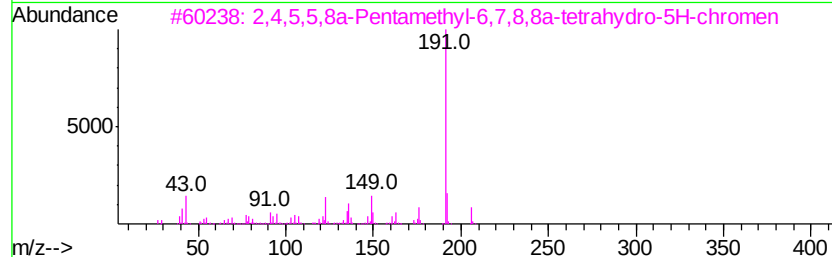
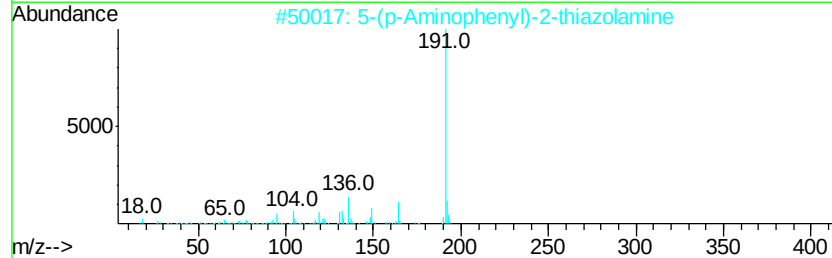
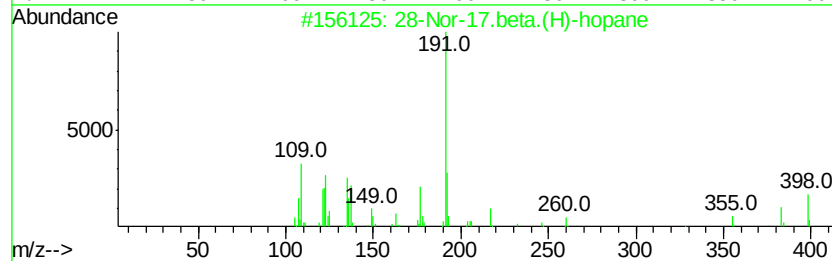
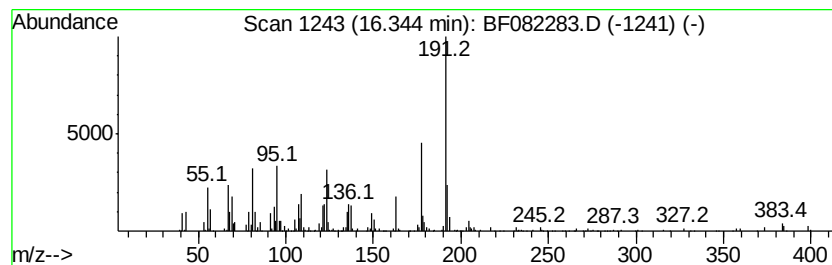
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 28-Nor-17.beta.(H)-hopane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.34	36.78 ng	713694	Perylene-d12	15.66

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	28-Nor-17.beta.(H)-hopane	398	C29H50	036728-72-0	64
2		5-(p-Aminophenyl)-2-thiazolamine	191	C9H9N3S	090349-87-4	38
3		2,4,5,5,8a-Pentamethyl-6,7,8,8a-...	206	C14H22O	1000195-40-9	38
4		.beta.-iso-Methyl ionone	206	C14H22O	1000285-40-2	38
5		1-Penten-3-one, 1-(2,6,6-trimeth...	206	C14H22O	000127-43-5	38



Data Path : Z:\HPCHEM1\BNA F\DATA\BF101515\
Data File : BF082283.D
Acq On : 14 Oct 2015 17:01
Operator : UM/IZ
Sample : G4015-08
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
LF-4

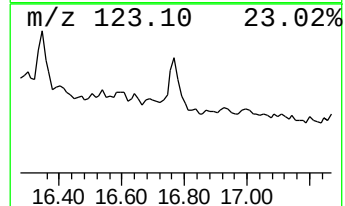
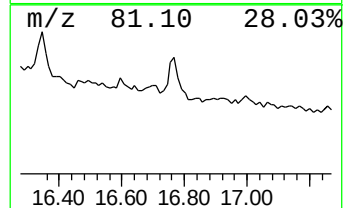
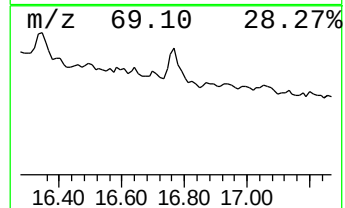
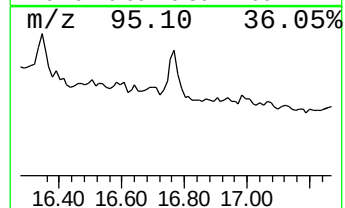
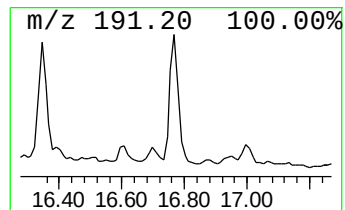
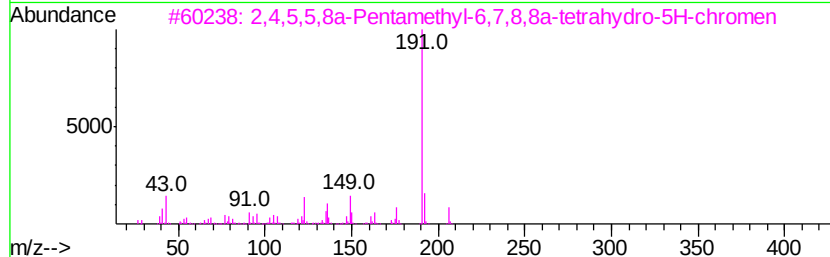
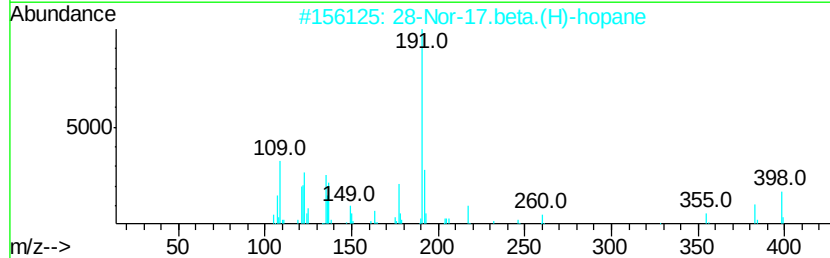
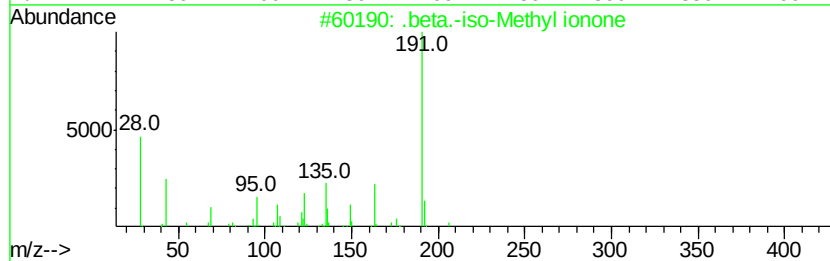
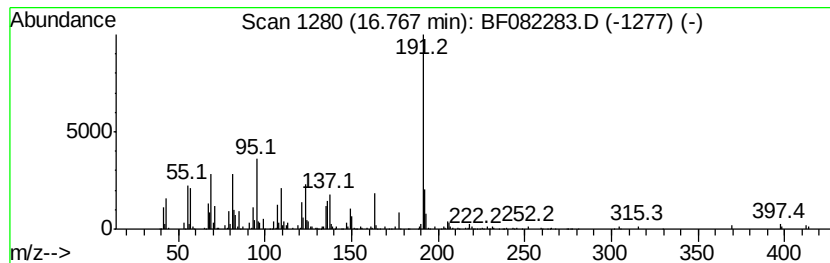
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF101015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 .beta.-iso-Methyl ionone Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.77	33.00 ng	640451	Perylene-d12	15.66

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	.beta.-iso-Methyl ionone	206	C14H22O	1000285-40-2	70
2		28-Nor-17.beta.(H)-hopane	398	C29H50	036728-72-0	50
3		2,4,5,5,8a-Pentamethyl-6,7,8,8a-...	206	C14H22O	1000195-40-9	50
4		1-Penten-3-one, 1-(2,6,6-trimeth...	206	C14H22O	000127-43-5	49
5		1-Cyclohexene, 1,3,3-trimethyl-2...	206	C14H22O	1000197-08-4	45



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF101515\
 Data File : BF082283.D
 Acq On : 14 Oct 2015 17:01
 Operator : UM/IZ
 Sample : G4015-08
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 LF-4

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
2-Pentanone, 4-hy...	4.99	59.9	ng	1184180	1	6.82	395727	20.0
unknown6.58	6.58	56.3	ng	1113130	1	6.82	395727	20.0
Hexadecane	10.77	3.8	ng	131803	4	11.36	697145	20.0
Dodecane, 1,1'-ox...	11.60	4.1	ng	142067	4	11.36	697145	20.0
Dodecane, 1-iodo-	11.65	3.6	ng	126183	4	11.36	697145	20.0
1-Decanol, 2-hexyl-	11.83	5.5	ng	192748	4	11.36	697145	20.0
Tritetracontane	11.91	4.6	ng	161127	4	11.36	697145	20.0
Eicosane	12.06	5.0	ng	172931	4	11.36	697145	20.0
Tetradecane, 2-me...	12.34	5.8	ng	202894	4	11.36	697145	20.0
unknown12.41	12.41	4.7	ng	164716	4	11.36	697145	20.0
Tetracosane	12.46	6.0	ng	211054	4	11.36	697145	20.0
Dichloroacetic ac...	12.64	3.1	ng	107253	4	11.36	697145	20.0
Nonadecane	12.66	6.4	ng	222414	4	11.36	697145	20.0
28-Nor-17.beta.(H...	16.34	36.8	ng	713694	6	15.66	388139	20.0
.beta.-iso-Methyl...	16.77	33.0	ng	640451	6	15.66	388139	20.0