

Data Path : Z:\HPCHEM1\BNA F\DATA\BF112215\
 Data File : BF083158.D
 Acq On : 22 Nov 2015 14:49
 Operator : UM/IZ
 Sample : G4504-06
 Misc :
 ALS Vial : 46 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW-05

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-BF112115.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	4.798	243	246	259	rBV	479410	888664	13.63%	1.598%
2	5.244	283	285	297	rBV	1943481	2510090	38.51%	4.515%
3	6.307	376	378	386	rBV	1414914	1704788	26.15%	3.066%
4	6.421	386	388	391	rBV	4774112	4597500	70.53%	8.269%
5	6.650	406	408	413	rVB	1297962	1247525	19.14%	2.244%
6	6.810	419	422	424	rBV	4252934	4670953	71.66%	8.401%
7	7.221	455	458	461	rBV	3794097	3894572	59.75%	7.005%
8	7.359	467	470	473	rVB	74455	86806	1.33%	0.156%
9	7.942	518	521	523	rBV	1460525	1611445	24.72%	2.898%
10	9.016	612	615	618	rBV	5094425	6518431	100.00%	11.724%
11	9.153	625	627	629	rBV	110829	103116	1.58%	0.185%
12	9.416	648	650	653	rBV	571351	537313	8.24%	0.966%
13	9.633	668	669	671	rVB	67623	66909	1.03%	0.120%
14	9.690	671	674	676	rBV	1451510	1848287	28.35%	3.324%
15	9.999	699	701	703	rBV2	39412	77197	1.18%	0.139%
16	10.136	709	713	715	rBV2	139674	242095	3.71%	0.435%
17	10.296	725	727	728	rBV2	70856	71290	1.09%	0.128%
18	10.479	740	743	745	rBV	3206707	3946490	60.54%	7.098%
19	10.605	752	754	758	rBV	156042	180085	2.76%	0.324%
20	10.868	775	777	778	rBV2	99240	105794	1.62%	0.190%
21	11.050	791	793	795	rBV2	134500	131029	2.01%	0.236%
22	11.165	801	803	805	rBV	1757479	1685857	25.86%	3.032%
23	11.302	814	815	817	rBV2	81644	78224	1.20%	0.141%
24	11.416	822	825	828	rVV	869999	883418	13.55%	1.589%
25	11.473	828	830	833	rVV2	144695	226507	3.47%	0.407%
26	11.725	850	852	855	rBV	1603305	1747623	26.81%	3.143%
27	12.228	894	896	899	rBV	3221277	3102567	47.60%	5.580%
28	12.273	899	900	903	rVB	89809	92464	1.42%	0.166%
29	12.422	910	913	915	rBV2	185402	331374	5.08%	0.596%
30	12.502	918	920	923	rBV	1130456	1298758	19.92%	2.336%
31	12.593	925	928	930	rBV	199815	248408	3.81%	0.447%
32	12.753	939	942	944	rBV	5500815	5779614	88.67%	10.395%
33	12.971	959	961	967	rVB2	117038	217972	3.34%	0.392%
34	13.211	980	982	984	rBV	356506	388488	5.96%	0.699%

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

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Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF112115.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	13.291	987	989	991	rBV2	190650	213617	3.28%	0.384%
36	13.656	1019	1021	1025	rVB	649662	680729	10.44%	1.224%
37	13.782	1030	1032	1035	rBV2	1090116	1564982	24.01%	2.815%
38	13.976	1047	1049	1054	rVB	84411	111656	1.71%	0.201%
39	14.216	1068	1070	1073	rVB	54757	68583	1.05%	0.123%
40	14.285	1074	1076	1085	rVB	87289	147708	2.27%	0.266%
41	14.571	1099	1101	1105	rVB	57007	78173	1.20%	0.141%
42	14.639	1105	1107	1109	rVB	152405	141451	2.17%	0.254%
43	14.868	1125	1127	1132	rVB	59040	83674	1.28%	0.150%
44	15.154	1149	1152	1155	rBV	988212	1175092	18.03%	2.114%
45	15.199	1155	1156	1165	rVB2	56014	122609	1.88%	0.221%
46	15.622	1191	1193	1196	rVB2	55374	89269	1.37%	0.161%

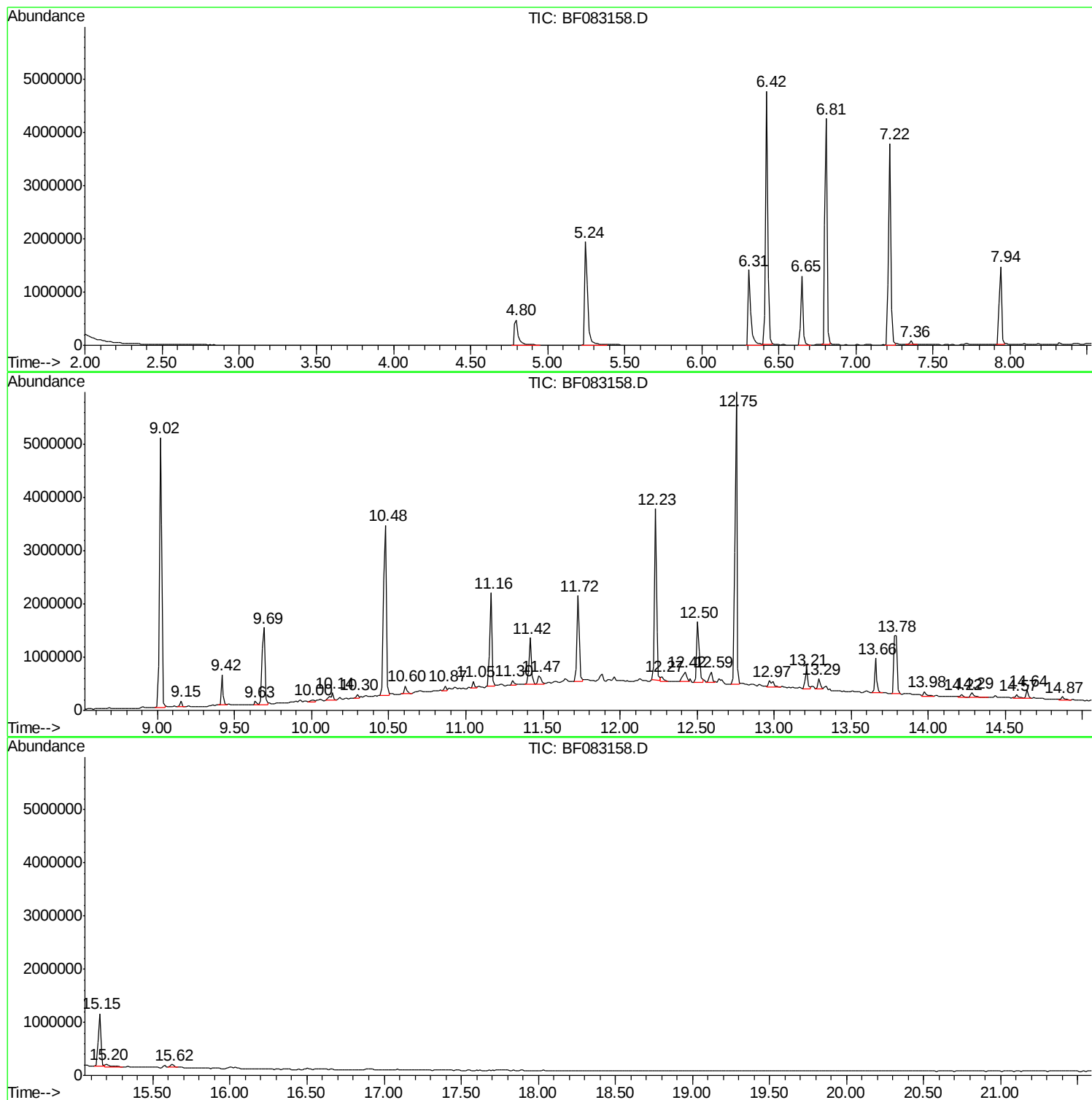
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF112115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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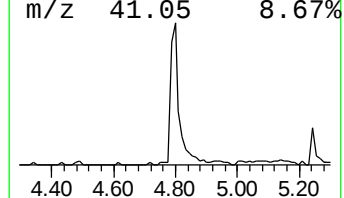
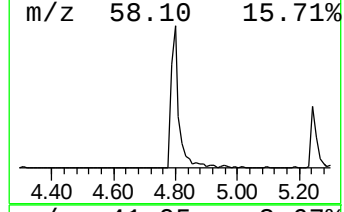
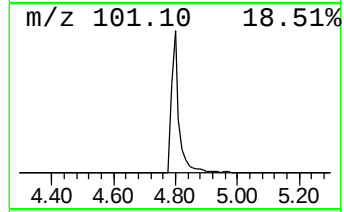
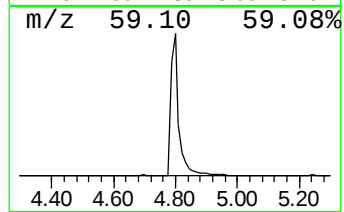
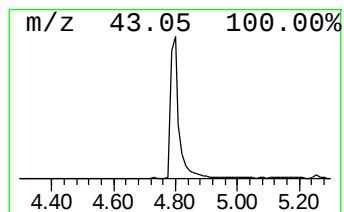
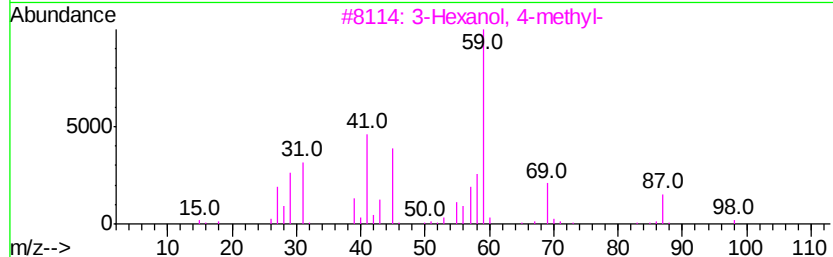
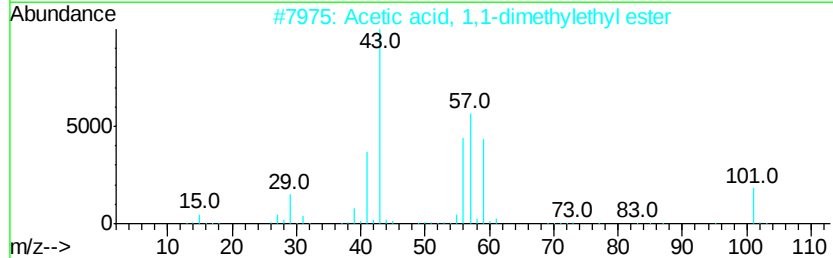
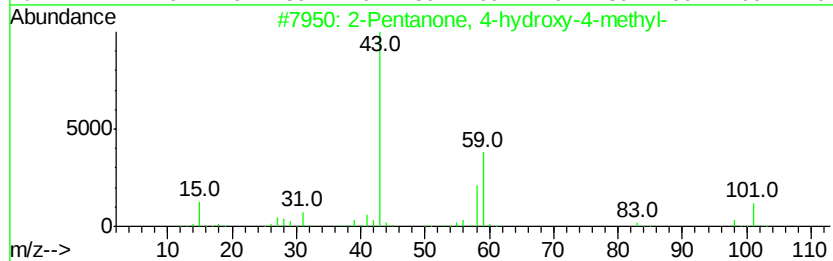
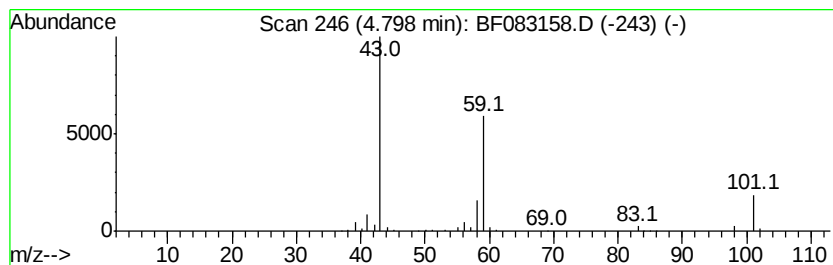
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF112115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.80	14.25 ng	888664	1,4-Dichlorobenzene-d4	6.65

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2		Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
3		3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
4		Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	23
5		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9



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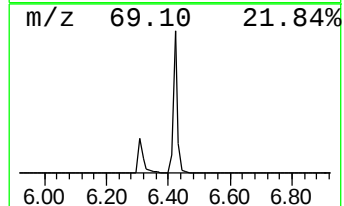
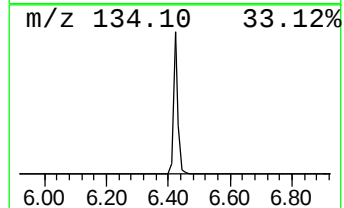
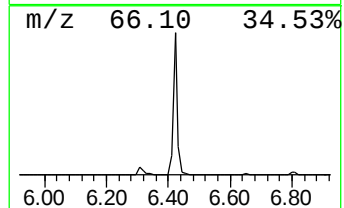
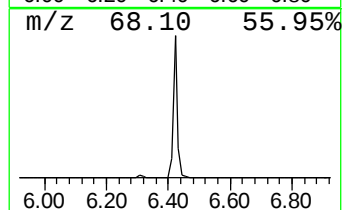
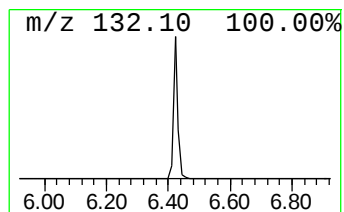
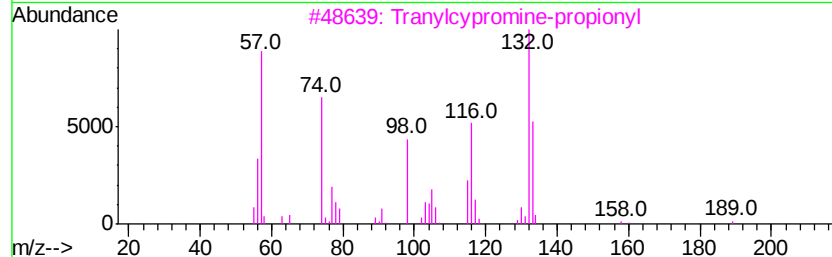
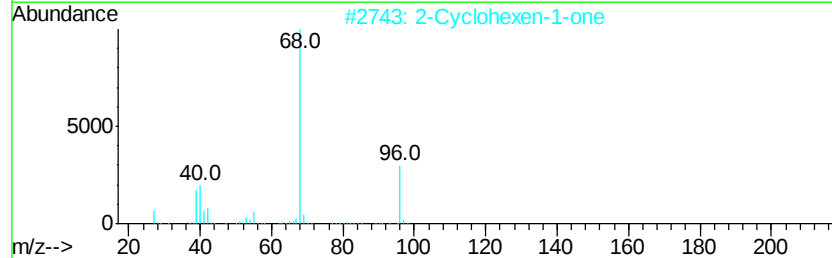
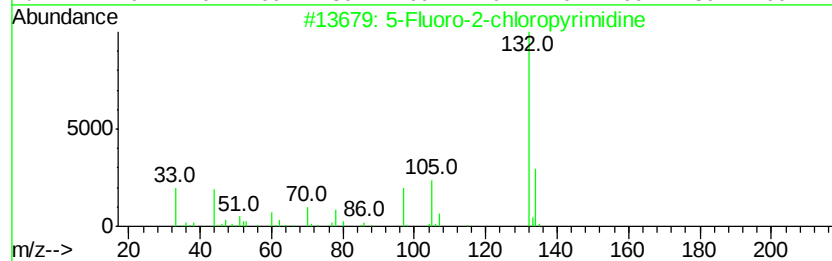
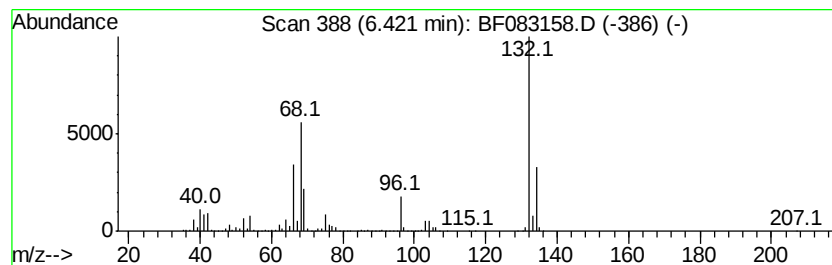
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 2 unknown6.42 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.42	73.71 ng	4597500	1,4-Dichlorobenzene-d4	6.65

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
2		2-Cyclohexen-1-one	96	C6H8O	000930-68-7	10
3		Tranvlcypromine-propionyl	189	C12H15NO	1000123-86-3	10
4		5-Aminoindole	132	C8H8N2	005192-03-0	10
5		(5-METHYL-2-PYRIDYL)ACETONITRILE	132	C8H8N2	1000241-93-9	10



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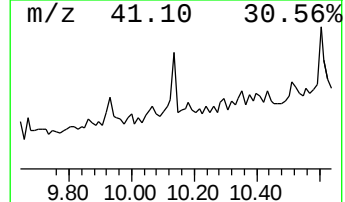
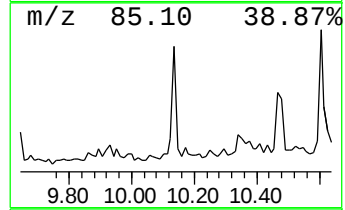
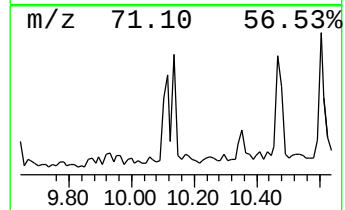
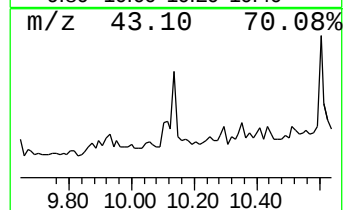
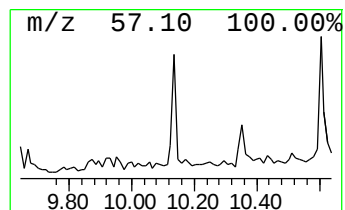
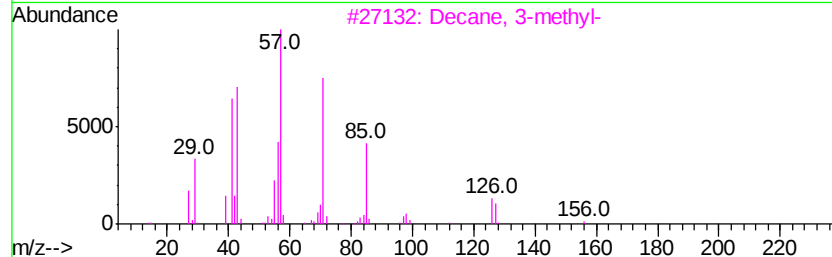
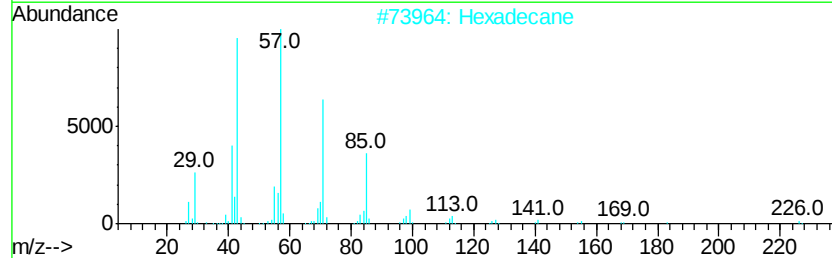
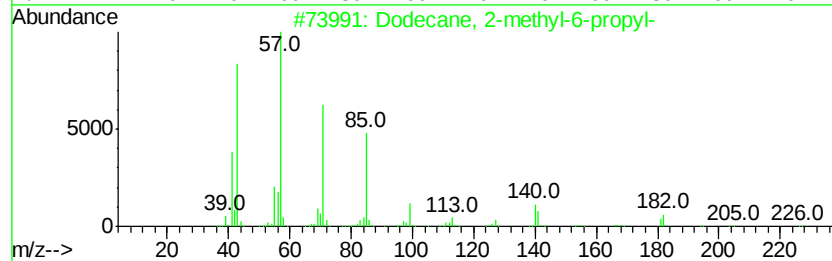
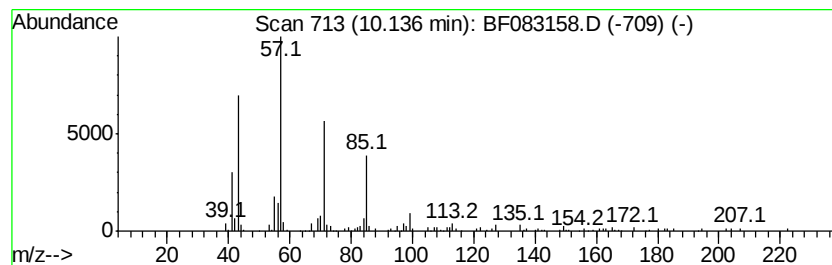
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 4 Dodecane, 2-methyl-6-propyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.14	2.62 ng	242095	Acenaphthene-d10	9.69
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Dodecane, 2-methyl-6-propyl-	226 C16H34	055045-08-4	86
2	Hexadecane	226 C16H34	000544-76-3	78
3	Decane, 3-methyl-	156 C11H24	013151-34-3	72
4	Decane, 2,3,8-trimethyl-	184 C13H28	062238-14-6	72
5	Heptacosane	380 C27H56	000593-49-7	72



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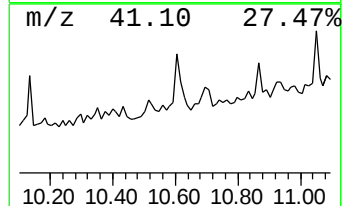
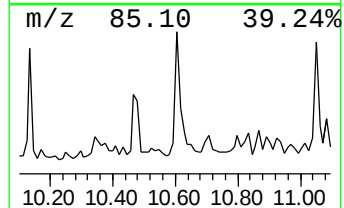
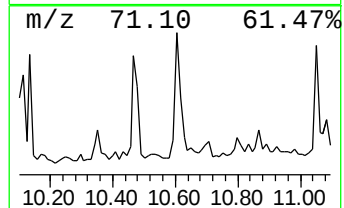
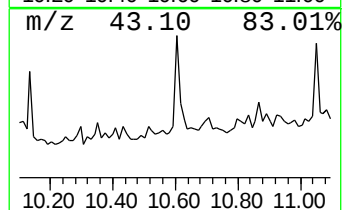
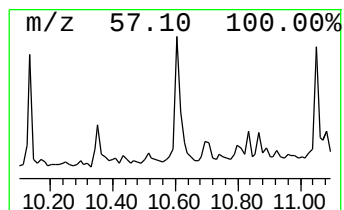
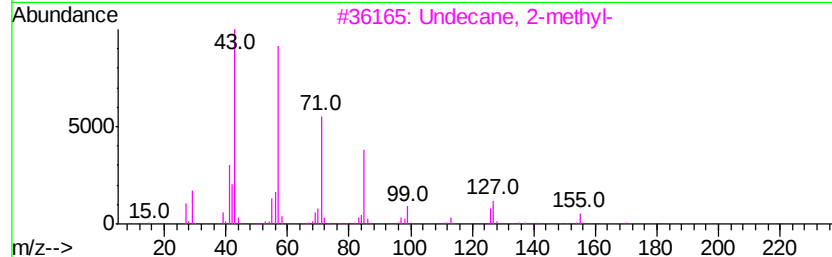
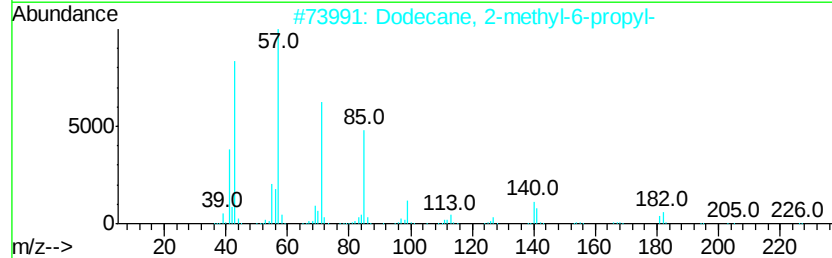
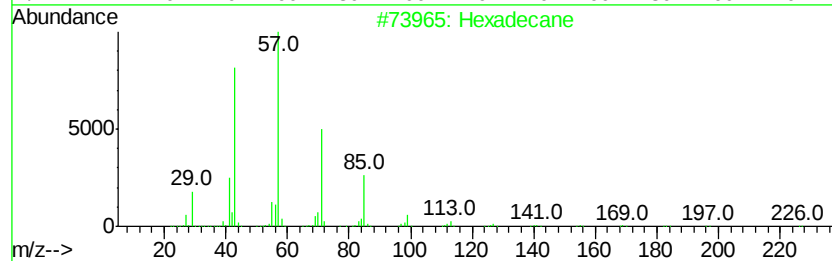
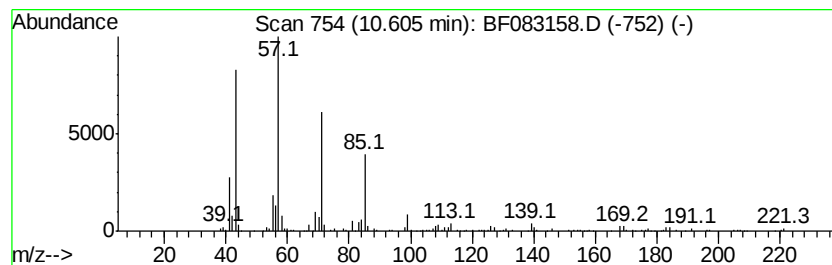
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Hexadecane Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.60	2.14 ng	180085	Phenanthrene-d10	11.16

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane		226	C16H34	000544-76-3	91
2	Dodecane, 2-methyl-6-propyl-		226	C16H34	055045-08-4	86
3	Undecane, 2-methyl-		170	C12H26	007045-71-8	74
4	Eicosane		282	C20H42	000112-95-8	72
5	Undecane, 3-methyl-		170	C12H26	001002-43-3	64



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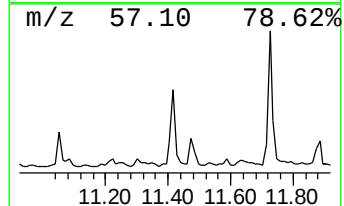
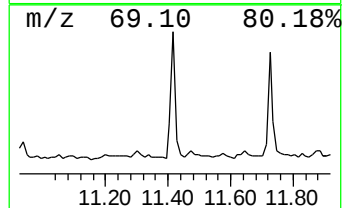
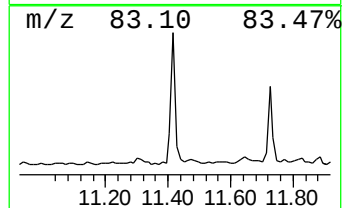
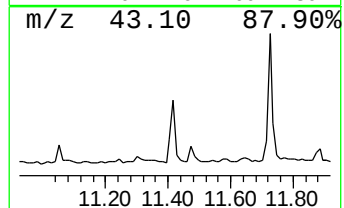
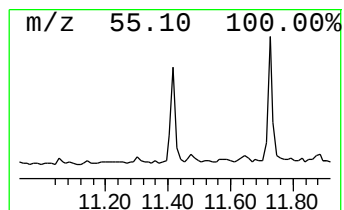
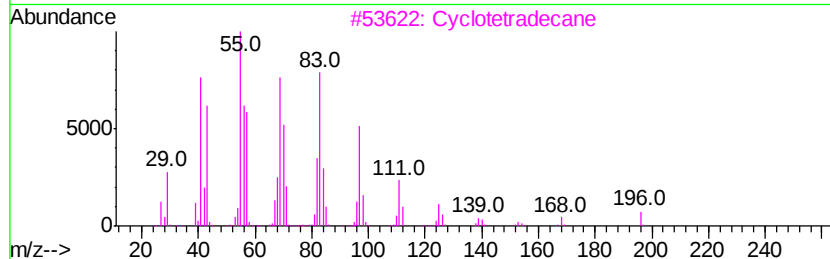
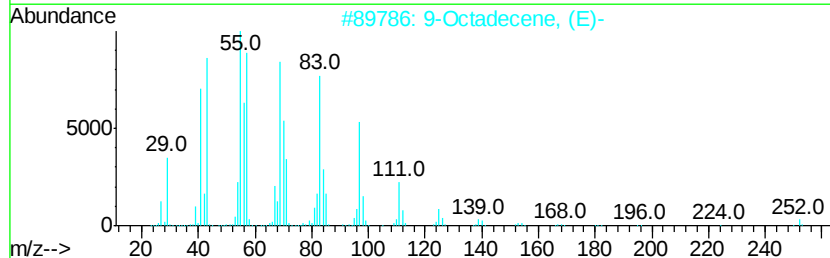
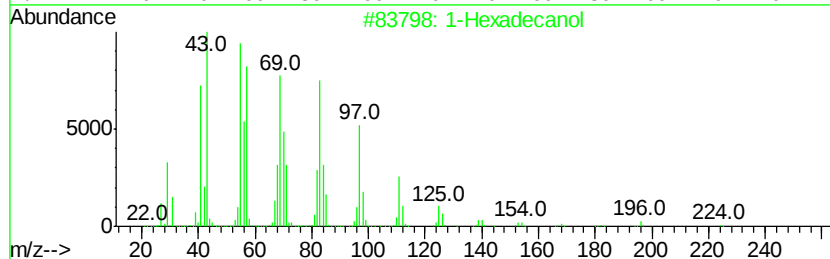
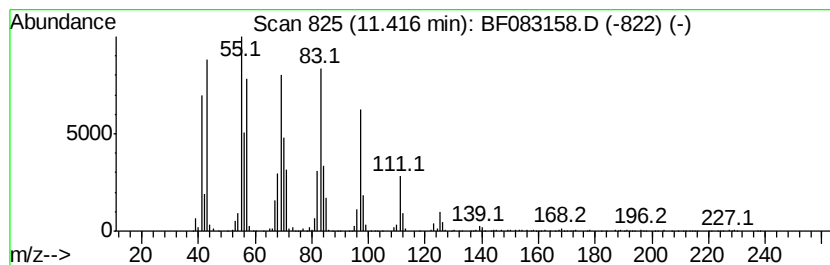
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 1-Hexadecanol Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.42	10.48 ng	883418	Phenanthrene-d10	11.16

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Hexadecanol	242	C16H34O	036653-82-4	94
2		9-Octadecene, (E)-	252	C18H36	007206-25-9	91
3		Cyclotetradecane	196	C14H28	000295-17-0	91
4		Cyclohexadecane	224	C16H32	000295-65-8	91
5		5-Octadecene, (E)-	252	C18H36	007206-21-5	91



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF112215\
Data File : BF083158.D
Acq On : 22 Nov 2015 14:49
Operator : UM/IZ
Sample : G4504-06
Misc :
ALS Vial : 46 Sample Multiplier: 1

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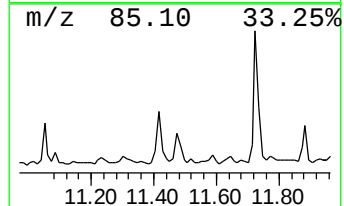
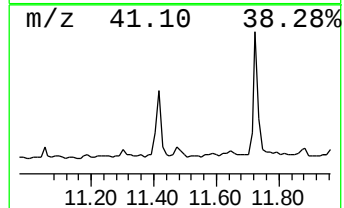
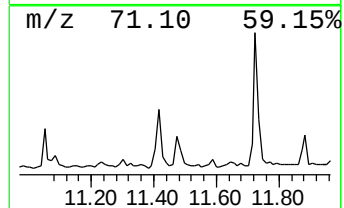
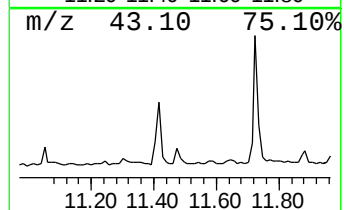
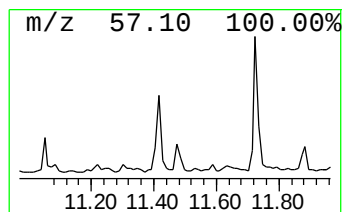
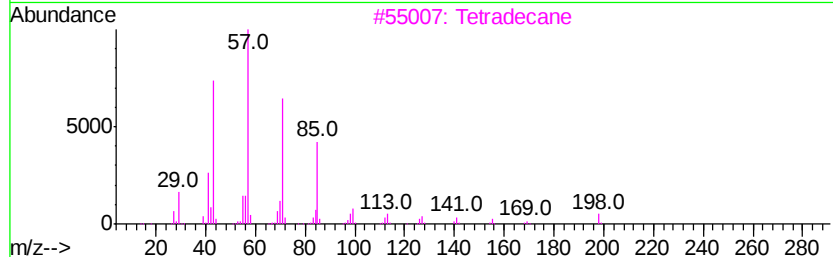
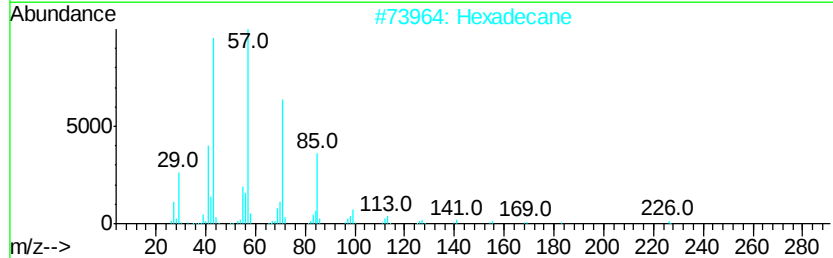
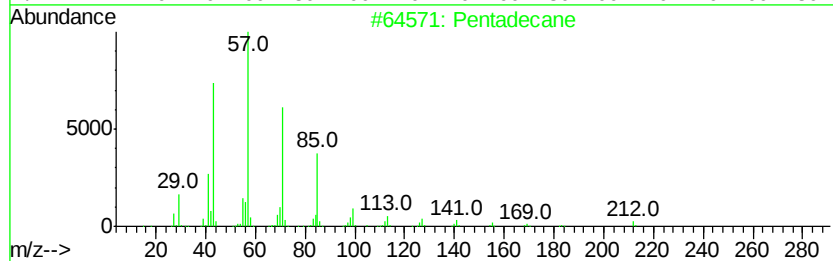
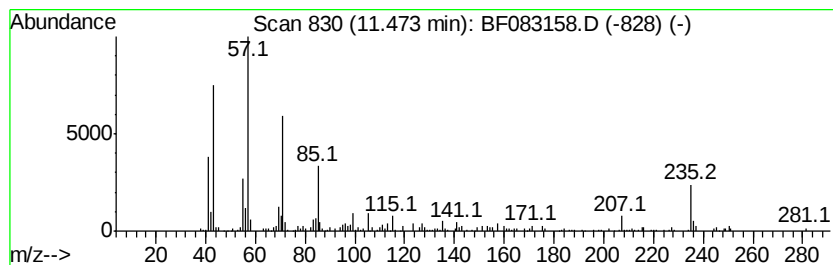
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Pentadecane Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.47	2.69 ng	226507	Phenanthrene-d10	11.16

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentadecane	212	C15H32	000629-62-9	93
2		Hexadecane	226	C16H34	000544-76-3	90
3		Tetradecane	198	C14H30	000629-59-4	70
4		Undecane, 2-methyl-	170	C12H26	007045-71-8	62
5		Tridecane, 1-iodo-	310	C13H27I	035599-77-0	58



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF112215\
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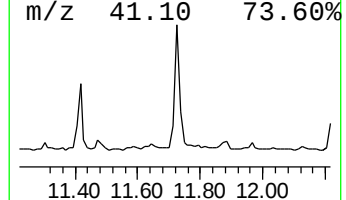
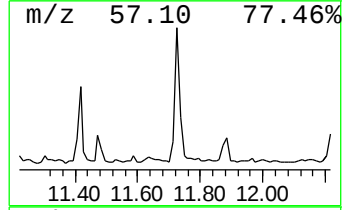
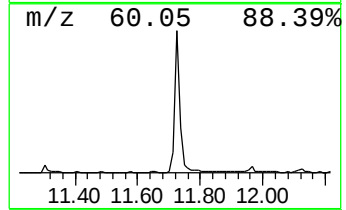
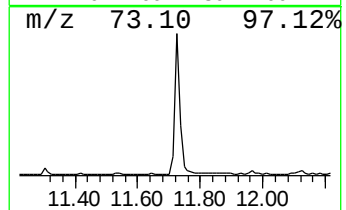
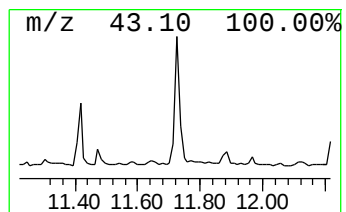
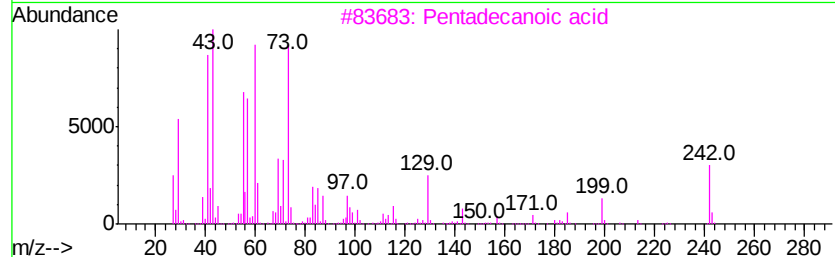
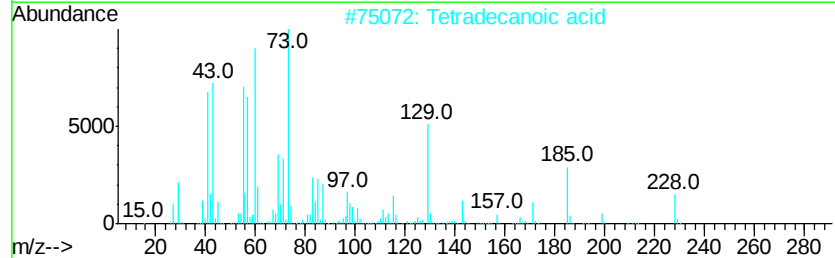
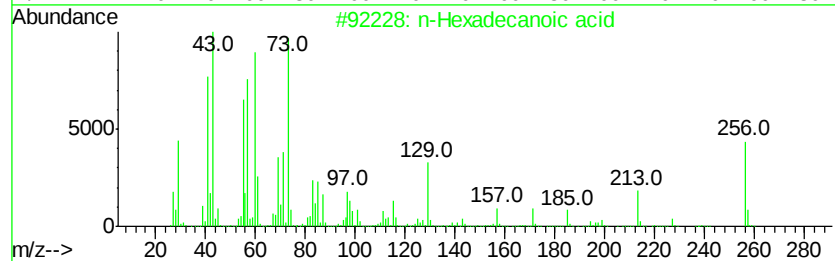
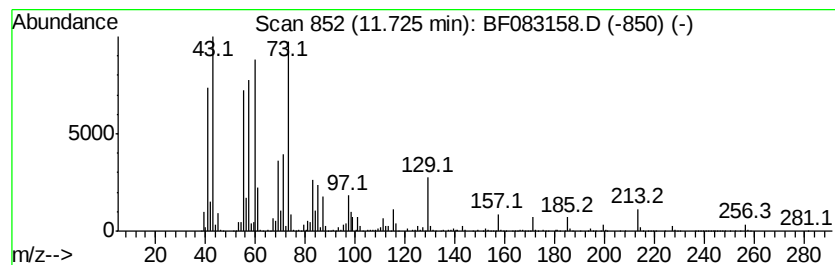
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 n-Hexadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.72	20.73 ng	1747620	Phenanthrene-d10	11.16

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2		Tetradecanoic acid	228	C14H28O2	000544-63-8	91
3		Pentadecanoic acid	242	C15H30O2	001002-84-2	90
4		Undecanoic acid	186	C11H22O2	000112-37-8	86
5		Tridecanoic acid	214	C13H26O2	000638-53-9	80



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF112215\
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Sample : G4504-06
Misc :
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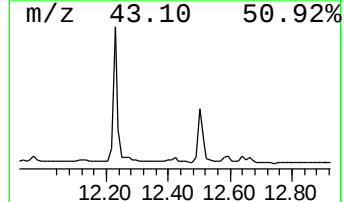
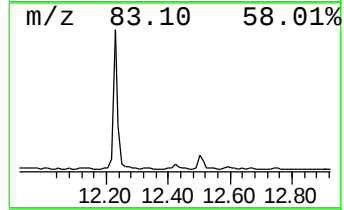
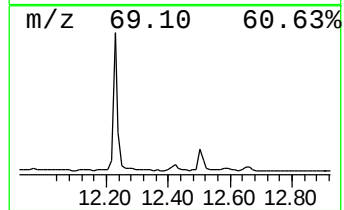
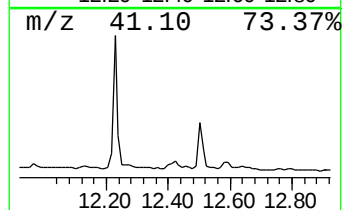
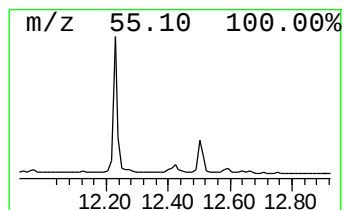
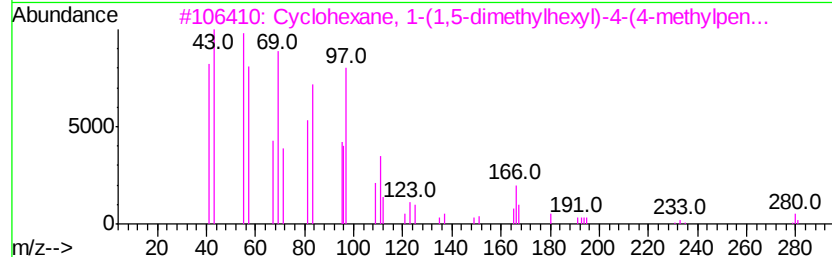
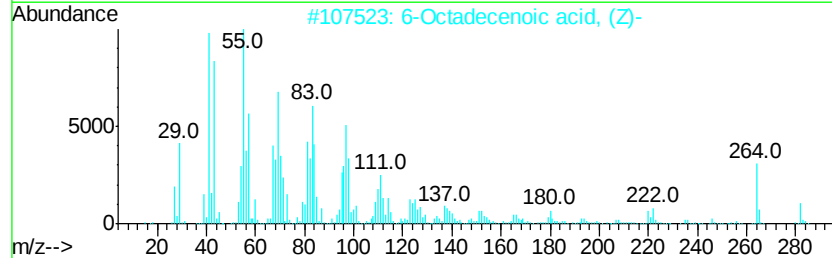
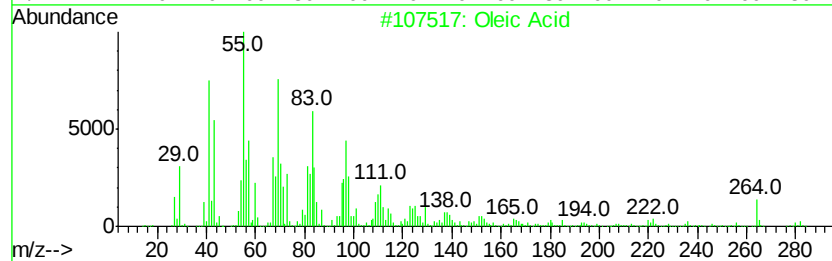
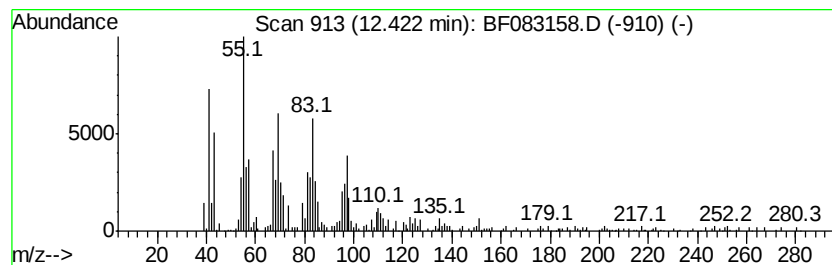
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Oleic Acid Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.42	3.93 ng	331374	Phenanthrene-d10	11.16	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Oleic Acid	282	C18H34O2	000112-80-1	86
2	6-Octadecenoic acid, (Z)-	282	C18H34O2	000593-39-5	60
3	Cyclohexane, 1-(1,5-dimethylhexyl)-4-(4-methylphenyl)-	280	C20H40	056009-20-2	55
4	2-Heptadecenal	252	C17H32O	1000143-48-6	53
5	E,Z-2,13-Octadecadien-1-ol	266	C18H34O	1000131-10-0	49



Data Path : Z:\HPCHEM1\BNA F\DATA\BF112215\
Data File : BF083158.D
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Misc :
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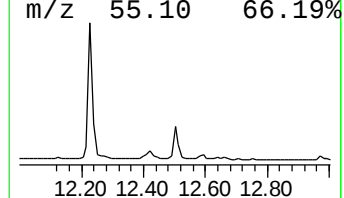
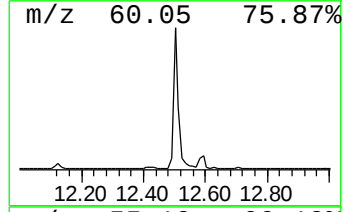
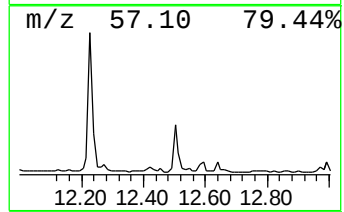
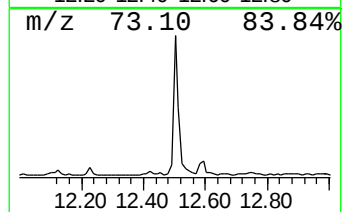
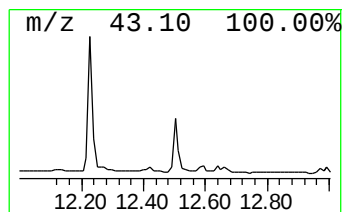
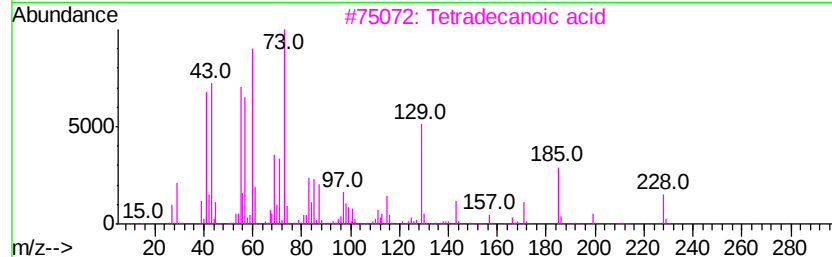
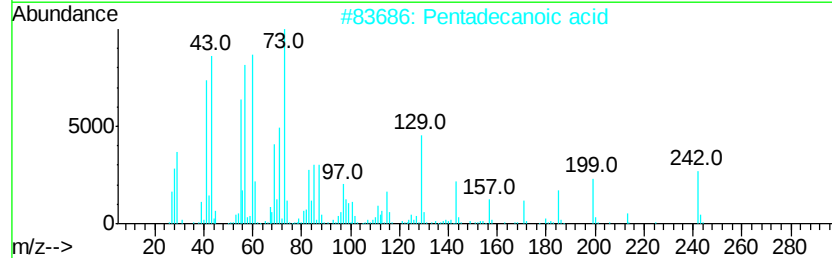
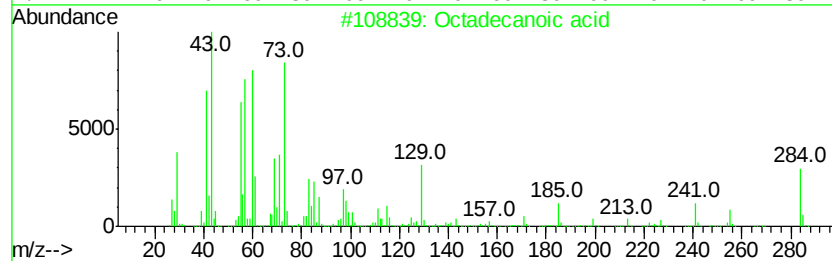
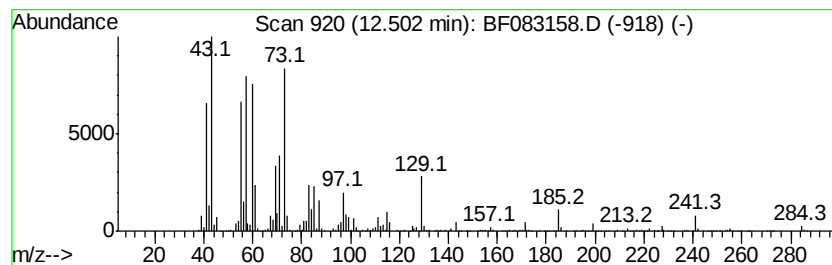
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 Octadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.50	16.60 ng	1298760	Chrysene-d12	13.78

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	95
2			Pentadecanoic acid	242	C15H30O2	001002-84-2	91
3			Tetradecanoic acid	228	C14H28O2	000544-63-8	62
4			Hydroxylamine, O-decyl-	173	C10H23NO	029812-79-1	20
5			Decane	142	C10H22	000124-18-5	11



Data Path : Z:\HPCHEM1\BNA F\DATA\BF112215\
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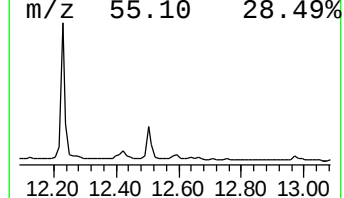
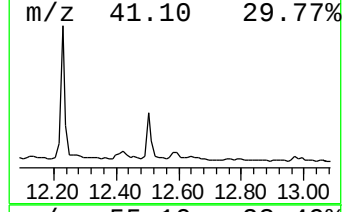
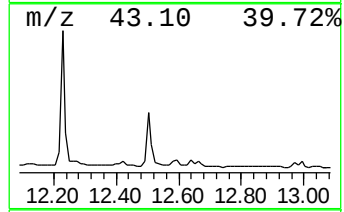
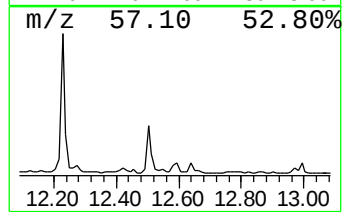
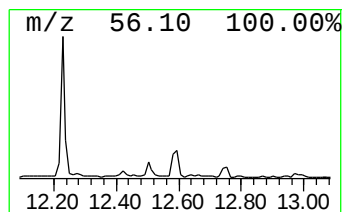
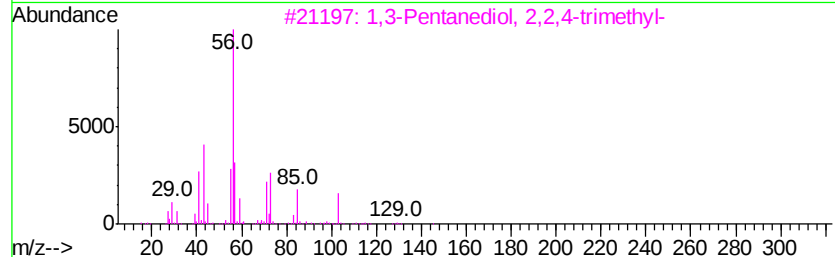
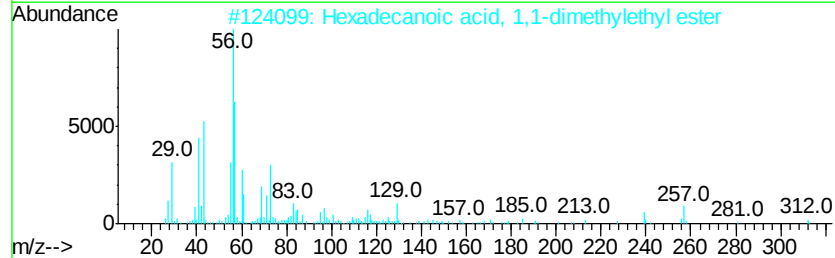
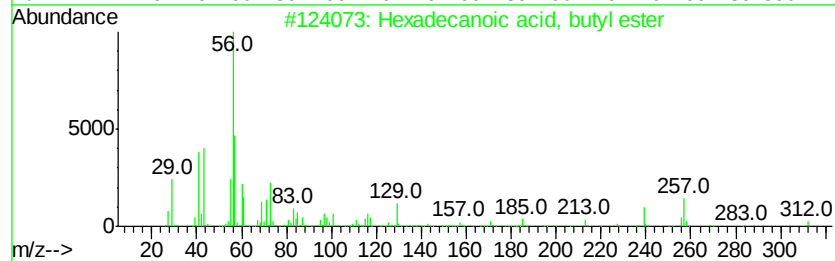
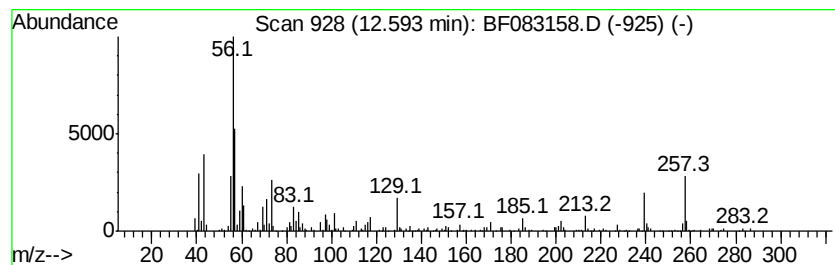
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Hexadecanoic acid, butyl ester Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.59	3.17 ng	248408	Chrysene-d12	13.78

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexadecanoic acid, butyl ester	312	C20H40O2	000111-06-8	72
2			Hexadecanoic acid, 1,1-dimethyle...	312	C20H40O2	031158-91-5	64
3			1,3-Pentanediol, 2,2,4-trimethyl-	146	C8H18O2	000144-19-4	43
4			Hexadecanoic acid, 2-methylpropyl...	312	C20H40O2	000110-34-9	37
5			Cyclopentanecarboxylic acid, 2-a...	129	C6H11NO2	037910-65-9	35



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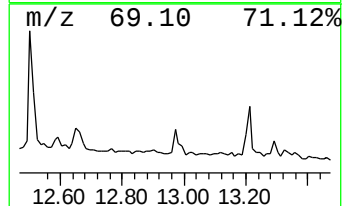
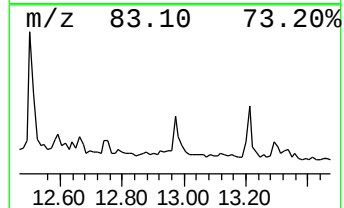
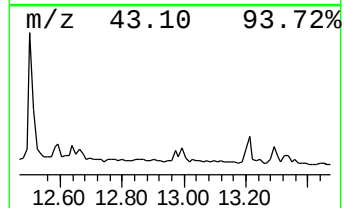
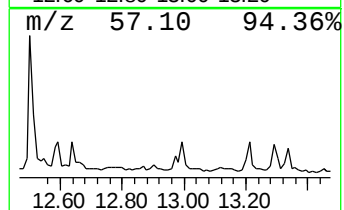
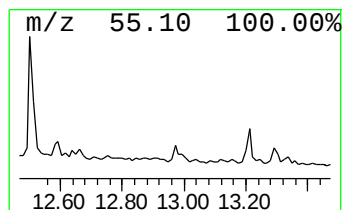
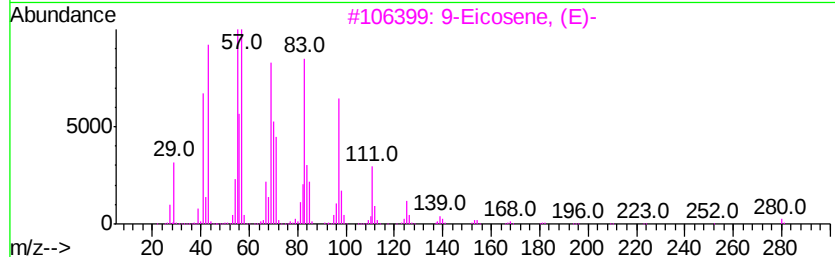
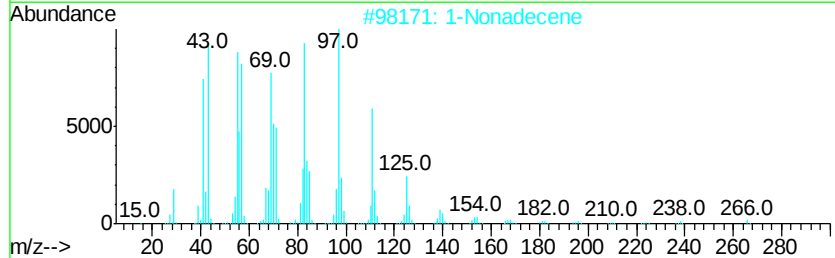
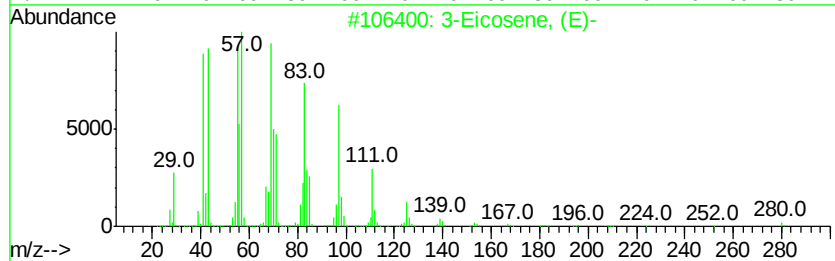
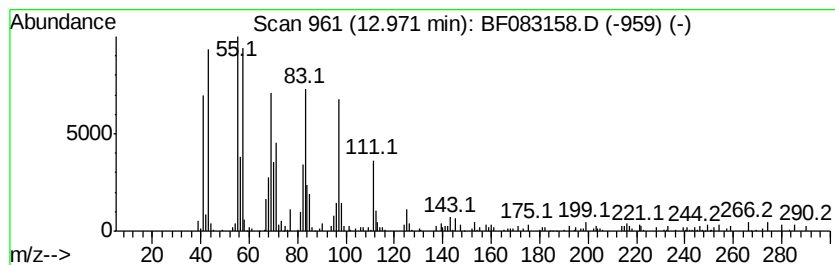
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 3-Eicosene, (E)- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.97	2.79 ng	217972	Chrysene-d12	13.78

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Eicosene, (E)-	280	C20H40	074685-33-9	99
2			1-Nonadecene	266	C19H38	018435-45-5	99
3			9-Eicosene, (E)-	280	C20H40	074685-29-3	96
4			1-Pentadecanethiol	244	C15H32S	025276-70-4	95
5			Z-5-Nonadecene	266	C19H38	1000131-11-8	95



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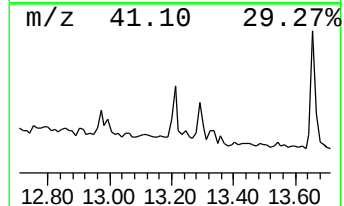
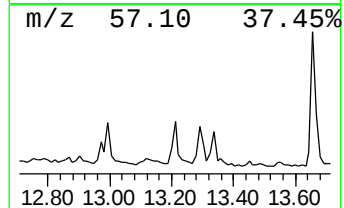
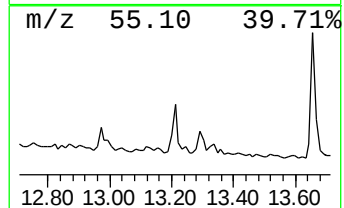
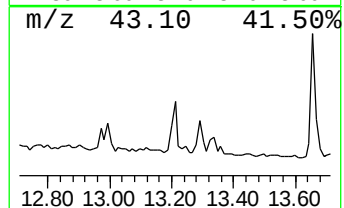
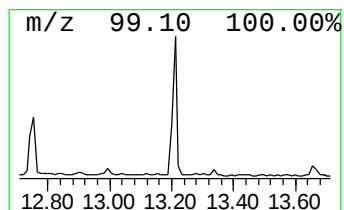
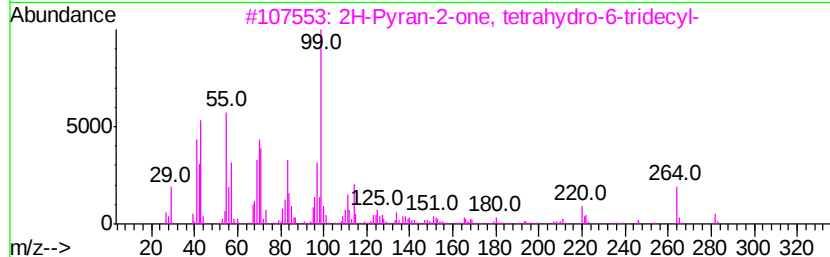
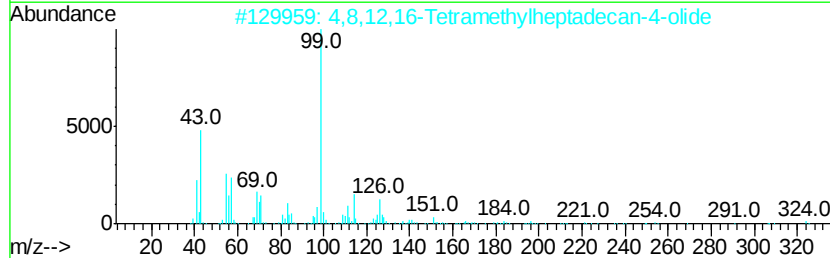
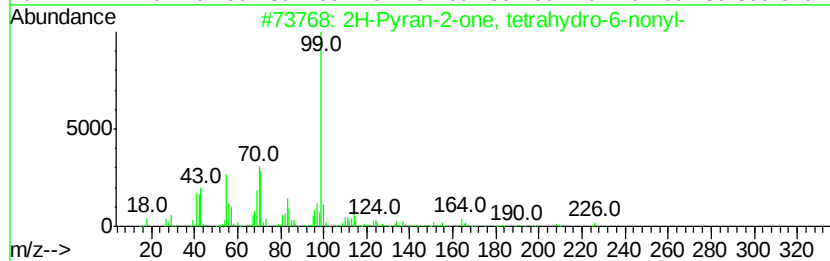
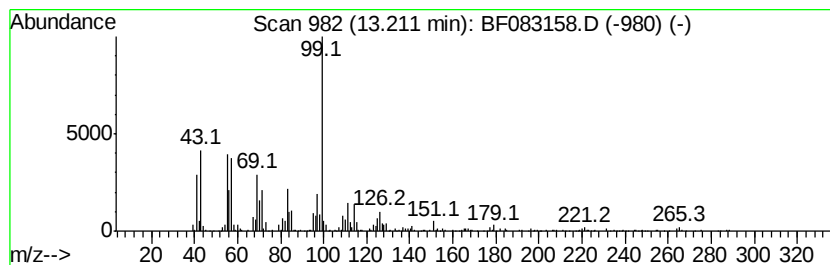
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 2H-Pyran-2-one, tetrahydro-... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.21	4.96 ng	388488	Chrysene-d12	13.78

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2H-Pyran-2-one, tetrahydro-6-nonyl-	226	C14H26O2	002721-22-4	68
2			4,8,12,16-Tetramethylheptadecan-...	324	C21H40O2	096168-15-9	58
3			2H-Pyran-2-one, tetrahydro-6-tri...	282	C18H34O2	001227-51-6	50
4			Spiro[1,3-dioxolane-2,2']-[6,7]di...	182	C9H14N2O2	1000142-34-9	47
5			.delta. Nonalactone	156	C9H16O2	003301-94-8	43



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF112215\
Data File : BF083158.D
Acq On : 22 Nov 2015 14:49
Operator : UM/IZ
Sample : G4504-06
Misc :
ALS Vial : 46 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MW-05

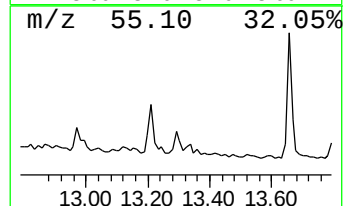
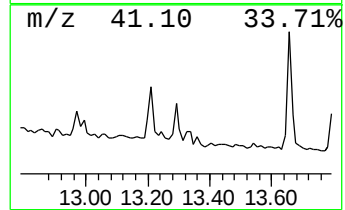
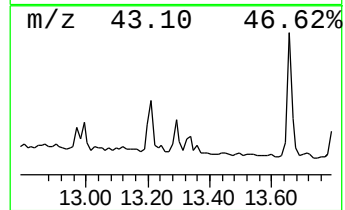
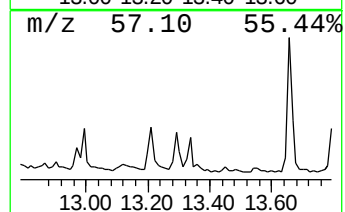
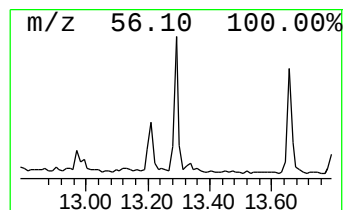
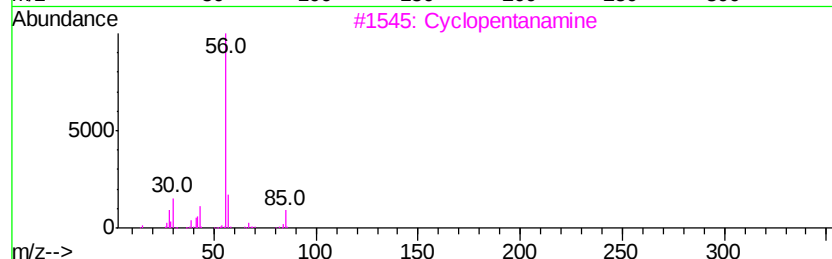
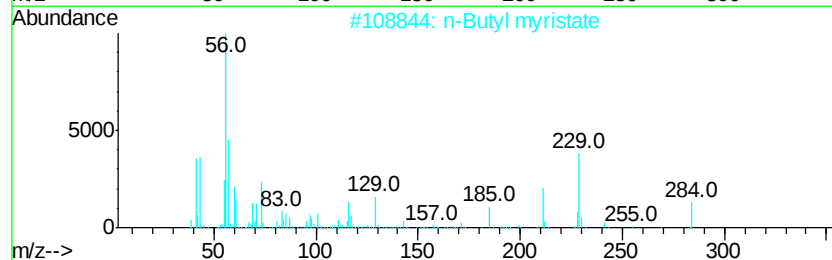
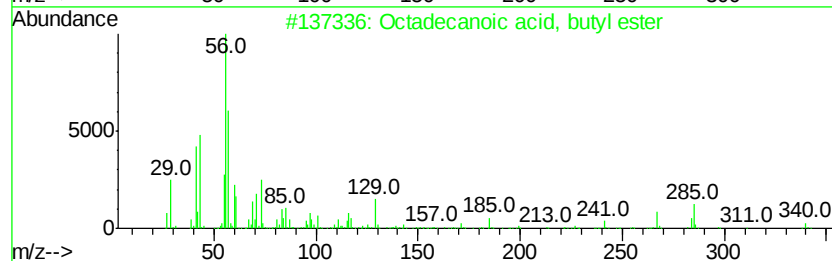
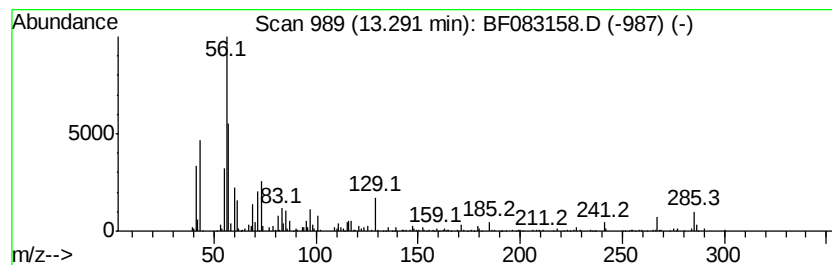
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Octadecanoic acid, butyl ester Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.29	2.73 ng	213617	Chrysene-d12	13.78

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecanoic acid, butyl ester	340	C22H44O2	000123-95-5	81
2		n-Butyl myristate	284	C18H36O2	000110-36-1	59
3		Cyclopentanamine	85	C5H11N	001003-03-8	35
4		Cyclobutane, 1-hexyl-2,3-dimethyl-	168	C12H24	055170-84-8	32
5		2,2-Dimethyl-1,3-butanediol	118	C6H14O2	000076-35-7	32



Data Path : Z:\HPCHEM1\BNA F\DATA\BF112215\
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Operator : UM/IZ
Sample : G4504-06
Misc :
ALS Vial : 46 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MW-05

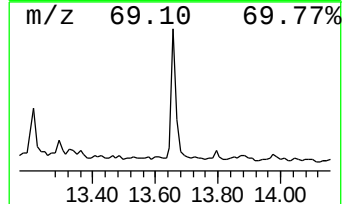
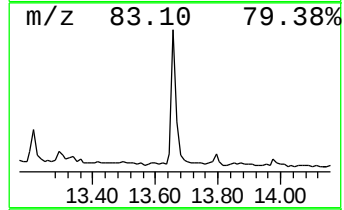
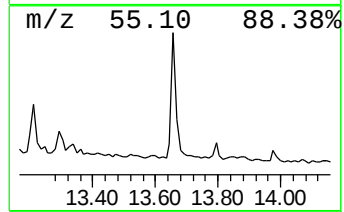
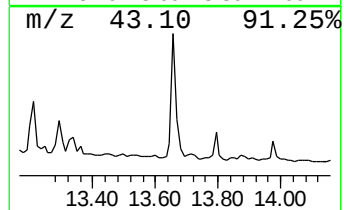
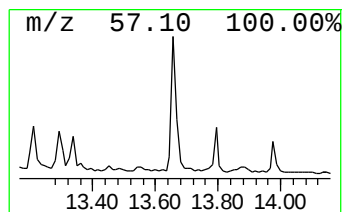
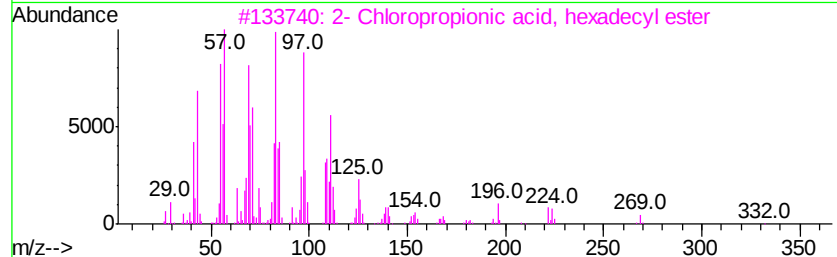
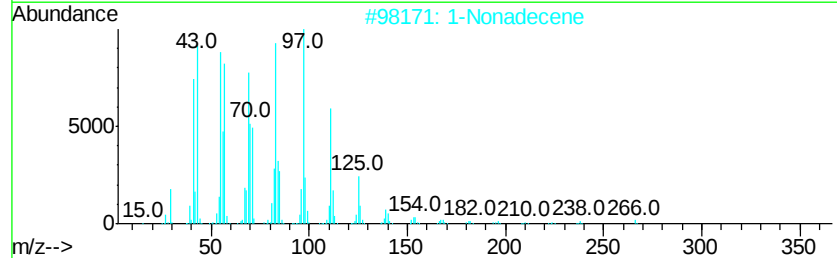
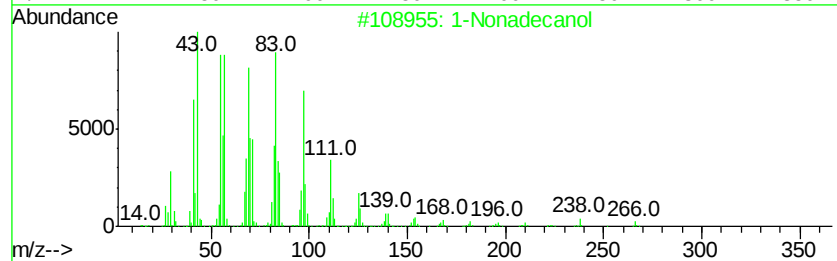
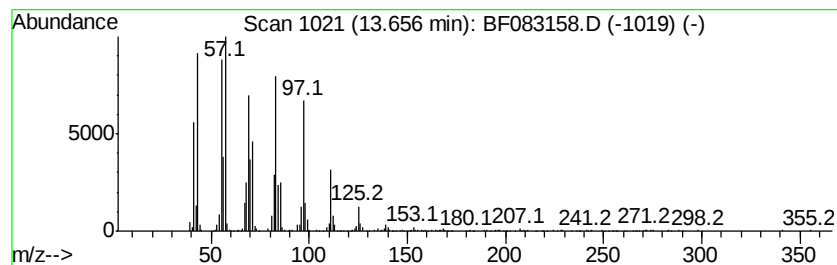
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 1-Nonadecanol Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.66	8.70 ng	680729	Chrysene-d12	13.78

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Nonadecanol	284	C19H40O	001454-84-8	94
2		1-Nonadecene	266	C19H38	018435-45-5	91
3		2- Chloropropionic acid, hexadec...	332	C19H37ClO2	086711-81-1	91
4		9-Eicosene, (E)-	280	C20H40	074685-29-3	91
5		Dichloroacetic acid, heptadecyl ...	366	C19H36Cl2O2	1000282-98-2	91



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Sample : G4504-06
Misc :
ALS Vial : 46 Sample Multiplier: 1

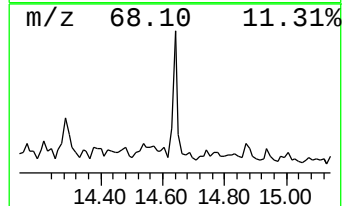
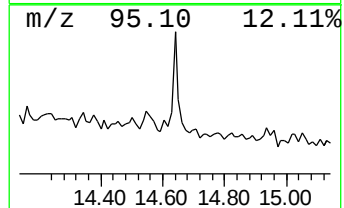
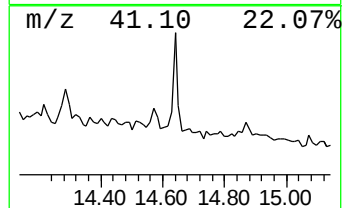
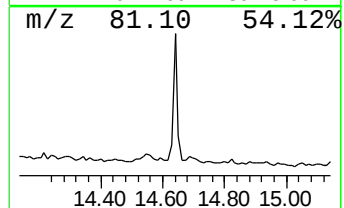
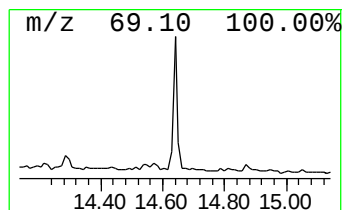
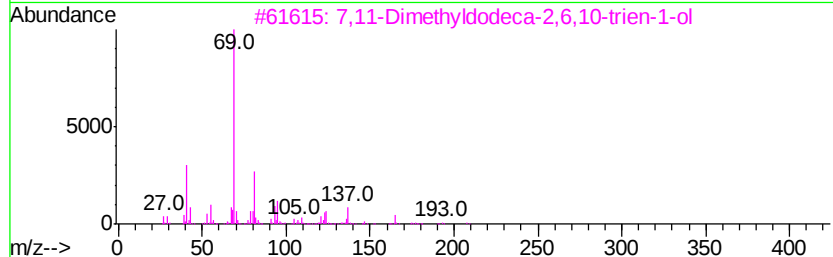
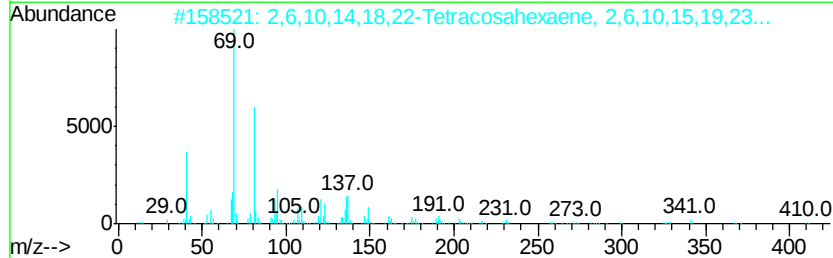
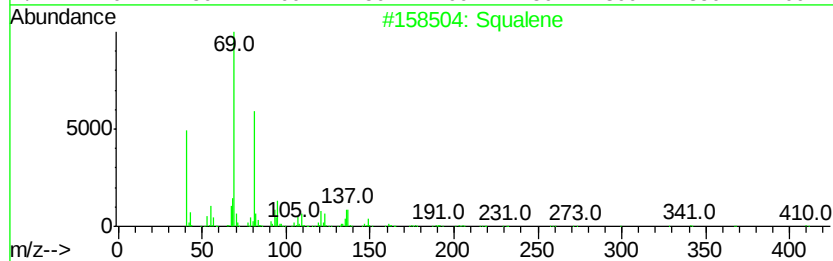
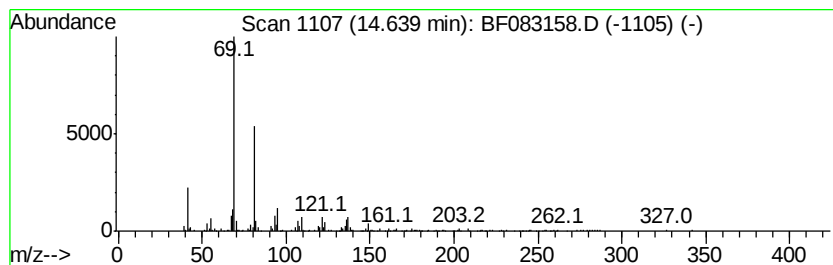
Instrument :
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF112115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 Squalene Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.64	2.41 ng	141451	Perylene-d12	15.15
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Squalene		410 C30H50	007683-64-9 91
2	2,6,10,14,18,22-Tetracosahexaene...		410 C30H50	000111-02-4 90
3	7,11-Dimethyldodeca-2,6,10-trien...		208 C14H240	125529-10-4 72
4	2,6,10-Dodecatrien-1-ol, 3,7,11-...		264 C17H2802	004128-17-0 64
5	1,5-Heptadiene, 3,3,6-trimethyl-		138 C10H18	035387-63-4 59



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF112215\
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 Sample : G4504-06
 Misc :
 ALS Vial : 46 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF112115.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...	4.80	14.3	ng	888664	1	6.65	1247530	20.0
unknown6.42	6.42	73.7	ng	4597500	1	6.65	1247530	20.0
Dodecane, 2-methy...	10.14	2.6	ng	242095	3	9.69	1848290	20.0
Hexadecane	10.60	2.1	ng	180085	4	11.16	1685860	20.0
1-Hexadecanol	11.42	10.5	ng	883418	4	11.16	1685860	20.0
Pentadecane	11.47	2.7	ng	226507	4	11.16	1685860	20.0
n-Hexadecanoic acid	11.72	20.7	ng	1747620	4	11.16	1685860	20.0
Oleic Acid	12.42	3.9	ng	331374	4	11.16	1685860	20.0
Octadecanoic acid	12.50	16.6	ng	1298760	5	13.78	1564980	20.0
Hexadecanoic acid...	12.59	3.2	ng	248408	5	13.78	1564980	20.0
3-Eicosene, (E)-	12.97	2.8	ng	217972	5	13.78	1564980	20.0
2H-Pyran-2-one, t...	13.21	5.0	ng	388488	5	13.78	1564980	20.0
Octadecanoic acid...	13.29	2.7	ng	213617	5	13.78	1564980	20.0
1-Nonadecanol	13.66	8.7	ng	680729	5	13.78	1564980	20.0
Squalene	14.64	2.4	ng	141451	6	15.15	1175090	20.0