

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF062416\
 Data File : BF088343.D
 Acq On : 24 Jun 2016 20:49
 Operator : UM/SJ
 Sample : H3602-15 10X
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 Q8-B3(0-8)C

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-BF060716.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	1.713	10	11	52	rVB	1047319	2600564	79.99%	20.694%
2	5.176	312	314	331	rBV3	55004	139077	4.28%	1.107%
3	6.250	407	408	416	rBV	69649	159146	4.90%	1.266%
4	6.353	416	417	435	rVV	123875	214464	6.60%	1.707%
5	6.582	435	437	449	rVV	372135	425723	13.09%	3.388%
6	6.742	449	451	460	rVB	98485	141710	4.36%	1.128%
7	7.165	486	488	501	rBV	50644	105609	3.25%	0.840%
8	7.873	548	550	552	rBV	482147	559391	17.21%	4.451%
9	8.948	642	644	651	rBV	234836	226617	6.97%	1.803%
10	9.348	678	679	685	rVB	40879	37312	1.15%	0.297%
11	9.611	700	702	705	rBV	597570	648202	19.94%	5.158%
12	10.399	770	771	778	rBB	69524	92443	2.84%	0.736%
13	11.085	829	831	841	rBV2	400509	564933	17.38%	4.495%
14	11.428	857	861	863	rBV	25265	46418	1.43%	0.369%
15	11.497	865	867	870	rBV	611223	710824	21.86%	5.656%
16	12.205	924	929	932	rBV2	1778000	3251155	100.00%	25.871%
17	12.285	934	936	939	rVB	317315	277333	8.53%	2.207%
18	12.674	968	970	972	rBV	130146	154120	4.74%	1.226%
19	12.914	989	991	995	rVB	34840	52728	1.62%	0.420%
20	13.577	1047	1049	1052	rVB	41397	48703	1.50%	0.388%
21	13.680	1052	1058	1059	rBV5	14011	35002	1.08%	0.279%
22	13.714	1059	1061	1064	rBV	513787	525352	16.16%	4.181%
23	13.897	1074	1077	1080	rBV	47943	65108	2.00%	0.518%
24	14.194	1101	1103	1109	rBV	51805	101621	3.13%	0.809%
25	14.297	1109	1112	1117	rBV5	22988	77492	2.38%	0.617%
26	14.491	1127	1129	1132	rBV	71641	90710	2.79%	0.722%
27	14.777	1152	1154	1156	rVB	53405	51655	1.59%	0.411%
28	15.074	1178	1180	1186	rBV	398372	541024	16.64%	4.305%
29	15.451	1211	1213	1215	rBV	42860	68787	2.12%	0.547%
30	15.623	1226	1228	1236	rVB2	61737	137111	4.22%	1.091%
31	15.863	1247	1249	1252	rVB	41977	65821	2.02%	0.524%
32	15.988	1257	1260	1265	rVB2	45173	109538	3.37%	0.872%
33	16.343	1288	1291	1295	rBV2	33613	71162	2.19%	0.566%
34	16.457	1299	1301	1303	rVV2	28099	45579	1.40%	0.363%

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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-BF060716.M
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35	16.514	1304	1306	1313	rVB4	31537	64314	1.98%	0.512%
36	17.566	1396	1398	1405	rVB4	18224	59971	1.84%	0.477%

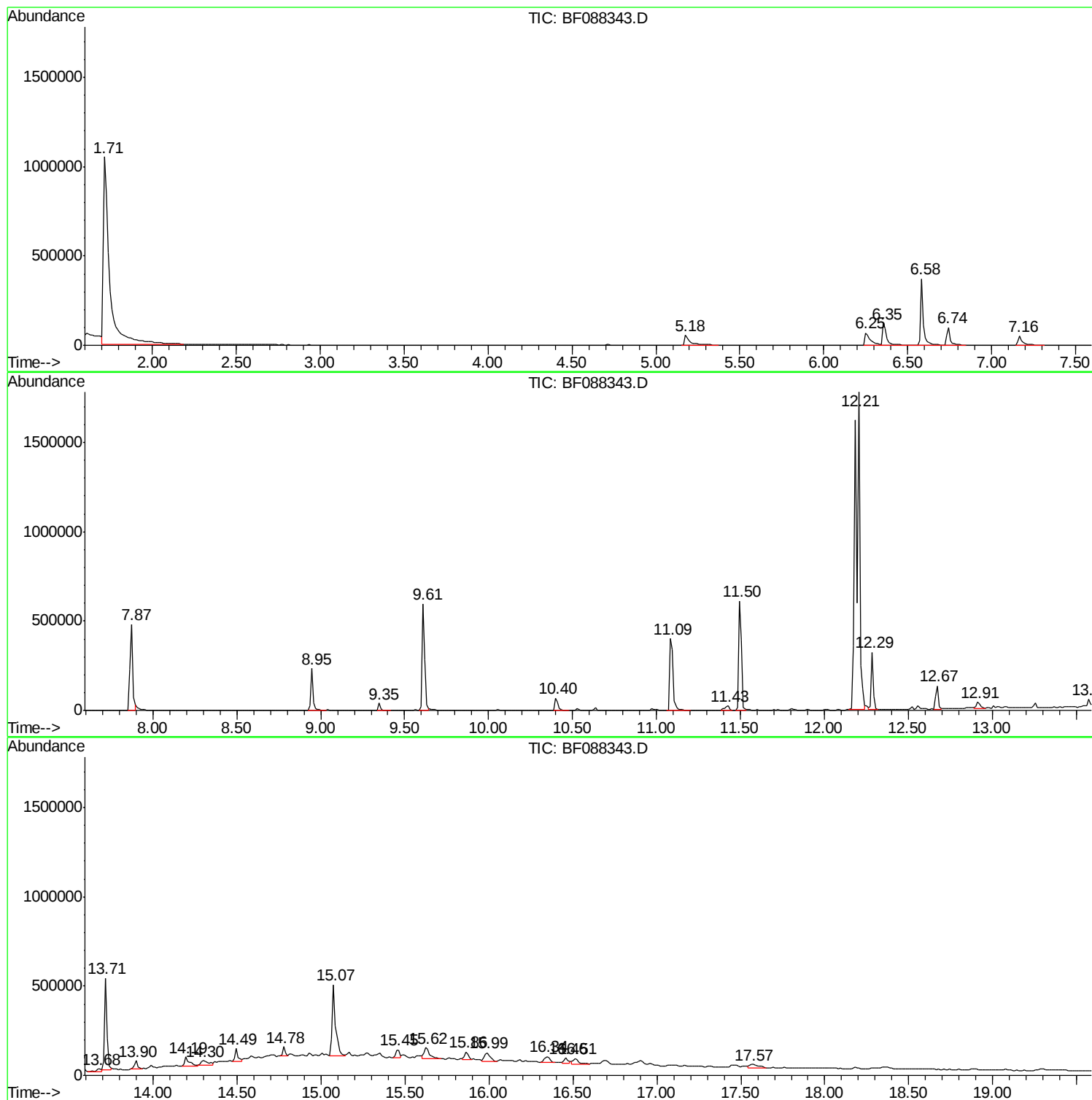
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060716.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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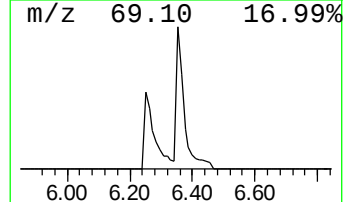
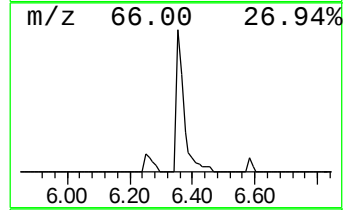
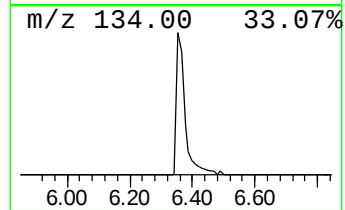
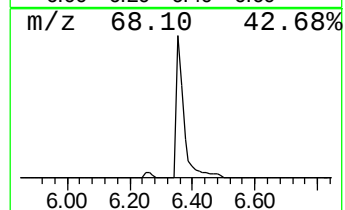
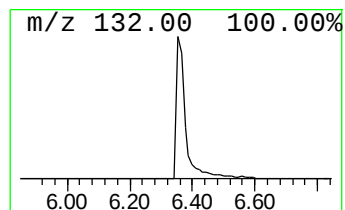
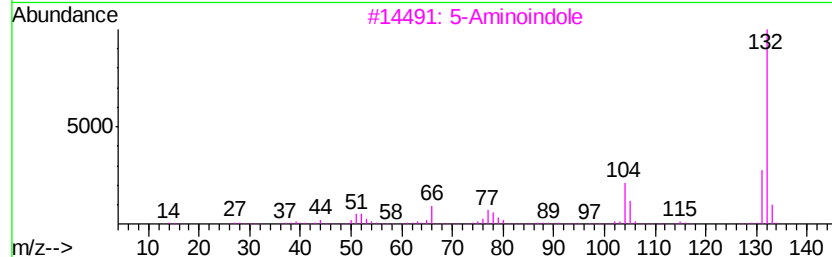
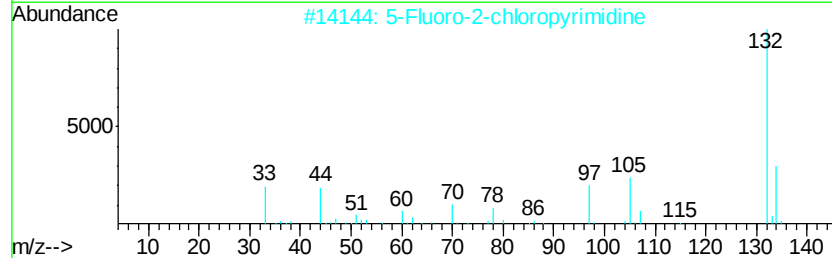
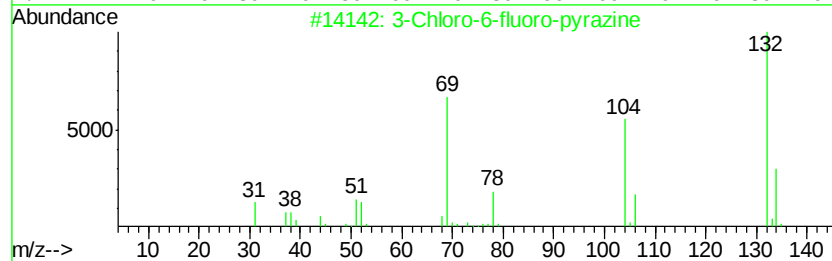
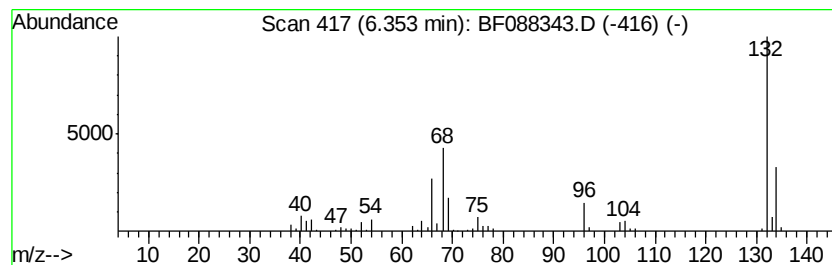
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown6.35 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.35	10.08 ng	214464	1,4-Dichlorobenzene-d4	6.58

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	10
4		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9
5		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9



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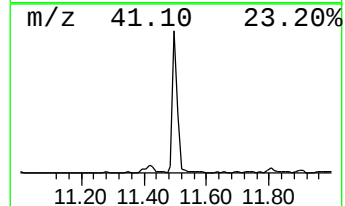
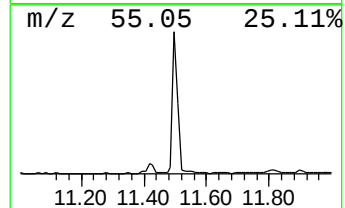
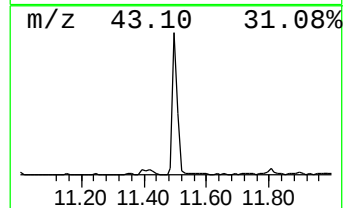
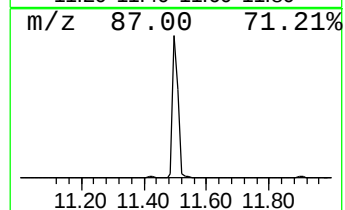
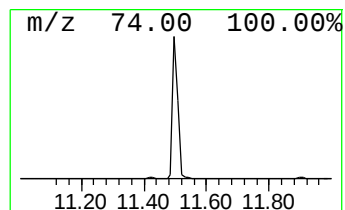
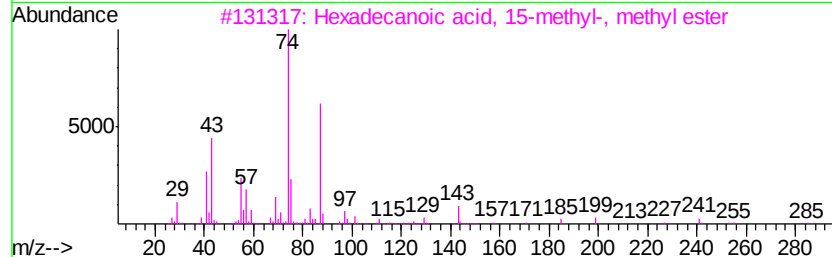
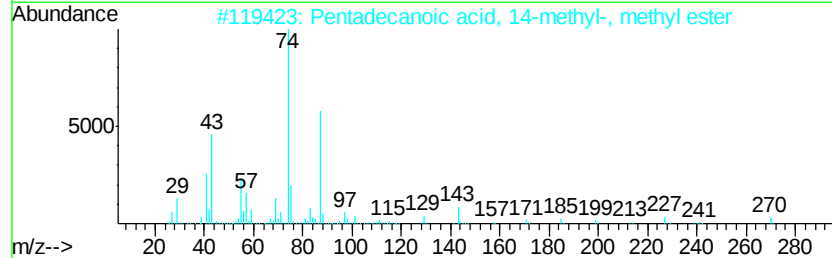
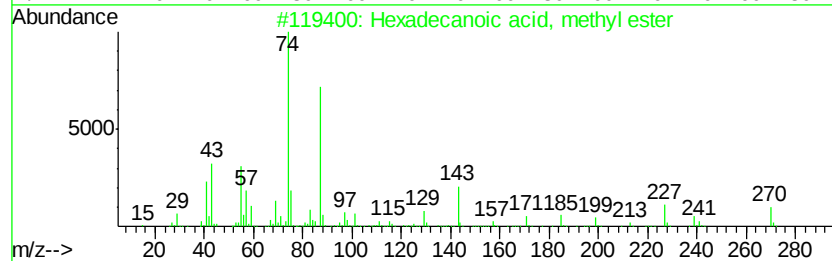
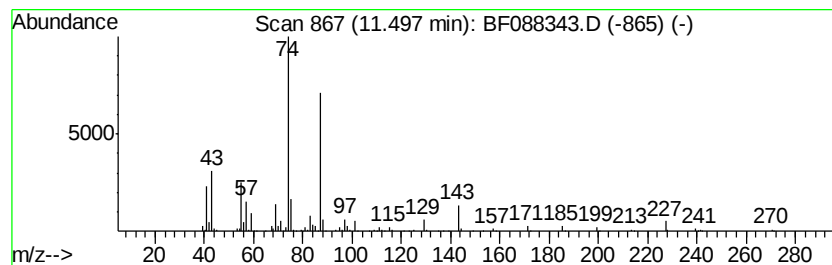
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TIC Integration Parameters: LSCINT.P

Peak Number 4 Hexadecanoic acid, methyl e... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.		
11.50	25.16 ng	710824	Phenanthrene-d10	11.09		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	98
2		Pentadecanoic acid, 14-methyl-, ...	270	C17H34O2	005129-60-2	93
3		Hexadecanoic acid, 15-methyl-, m...	284	C18H36O2	006929-04-0	91
4		Methyl tetradecanoate	242	C15H30O2	000124-10-7	90
5		Methyl 11-methyl-dodecanoate	228	C14H28O2	1000336-45-1	90



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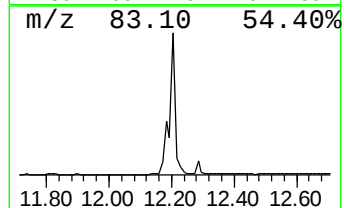
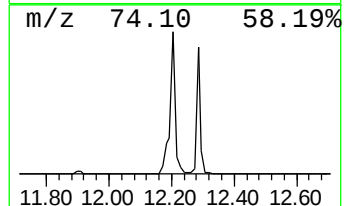
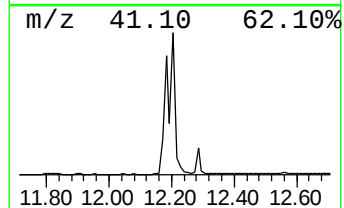
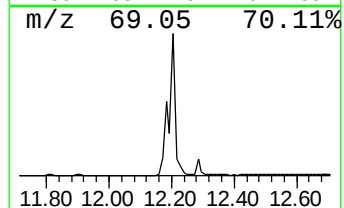
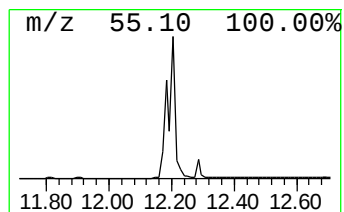
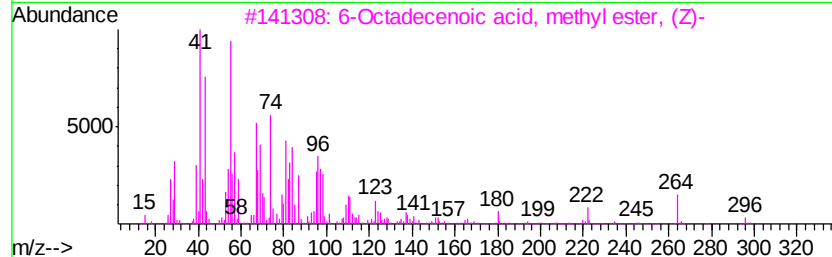
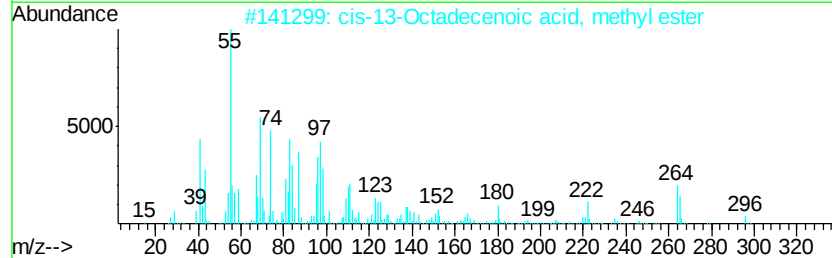
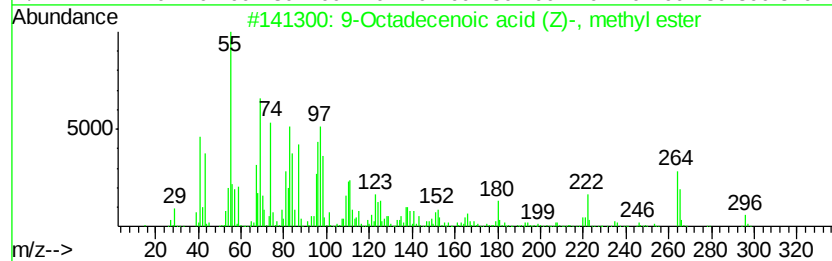
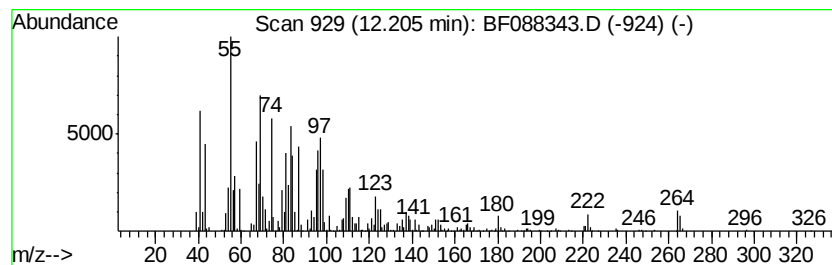
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 9-Octadecenoic acid (Z)-, m... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.21	115.10 ng	3251160	Phenanthrene-d10	11.09

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		9-Octadecenoic acid (Z)-, methyl...	296	C19H36O2	000112-62-9	99
2		cis-13-Octadecenoic acid, methyl...	296	C19H36O2	1000333-58-3	99
3		6-Octadecenoic acid, methyl este...	296	C19H36O2	002777-58-4	99
4		trans-13-Octadecenoic acid, meth...	296	C19H36O2	1000333-61-3	99
5		8-Octadecenoic acid, methyl ester	296	C19H36O2	002345-29-1	99



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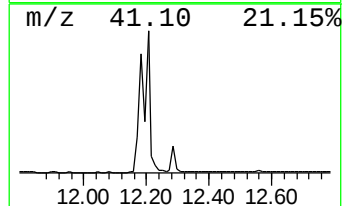
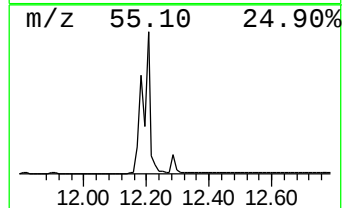
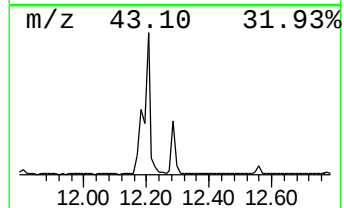
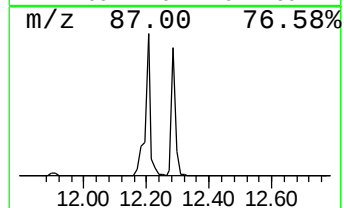
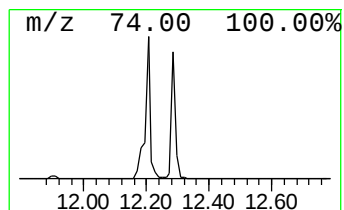
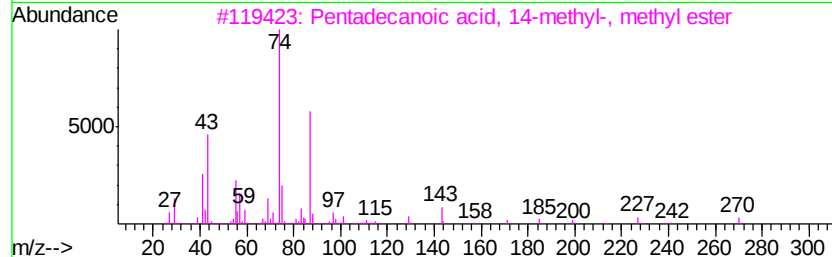
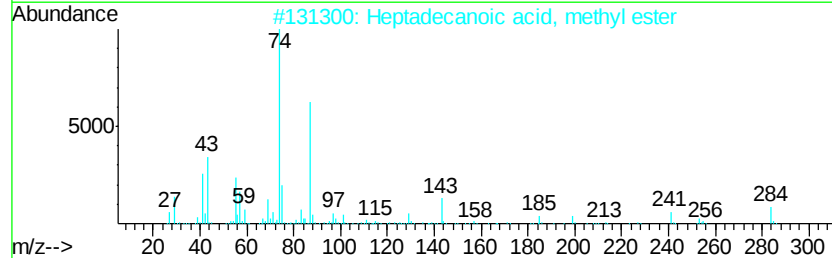
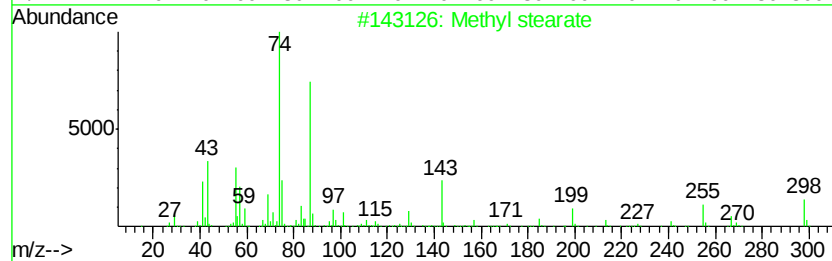
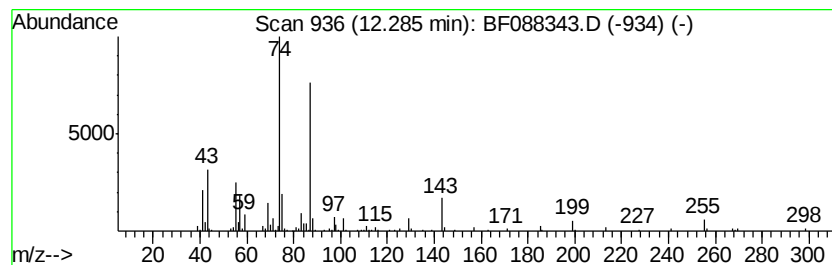
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TIC Library : C:\Database\NIST11.L
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Peak Number 6 Methyl stearate Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.29	9.82 ng	277333	Phenanthrene-d10	11.09	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Methyl stearate	298	C19H38O2	000112-61-8	99
2	Heptadecanoic acid, methyl ester	284	C18H36O2	001731-92-6	91
3	Pentadecanoic acid, 14-methyl-, ...	270	C17H34O2	005129-60-2	91
4	Methyl tetradecanoate	242	C15H30O2	000124-10-7	91
5	Hexadecanoic acid, 15-methyl-, m...	284	C18H36O2	006929-04-0	91



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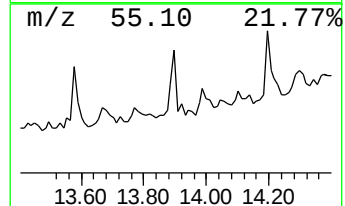
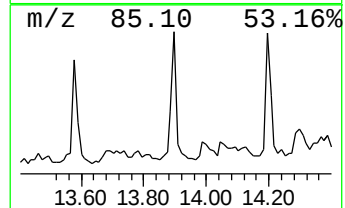
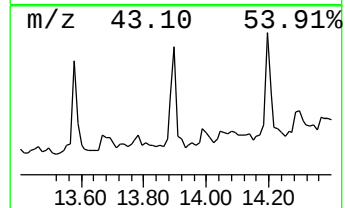
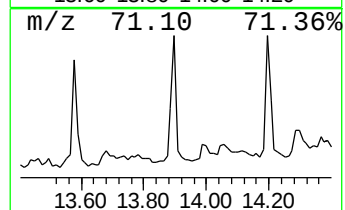
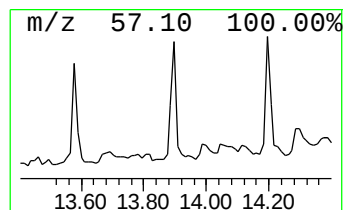
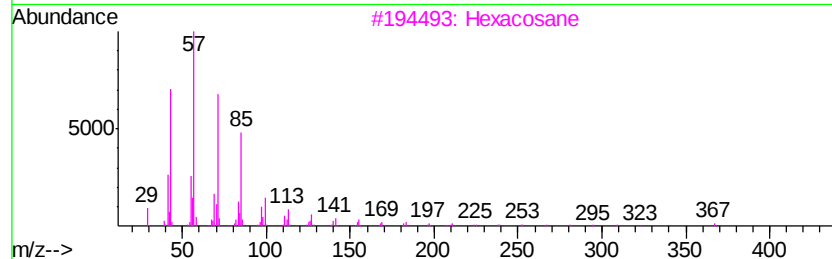
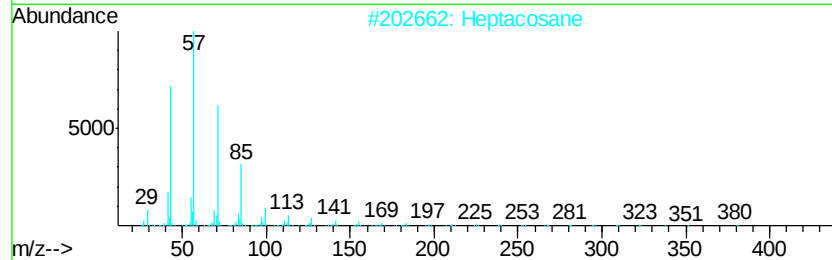
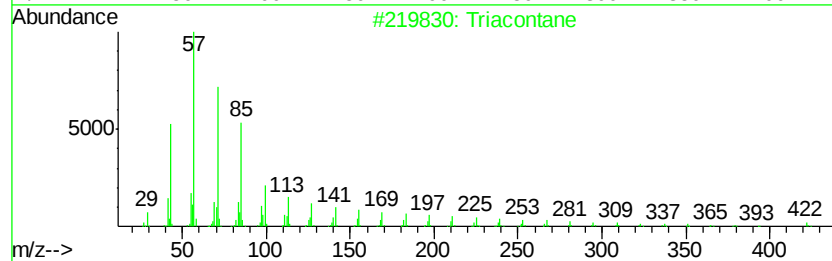
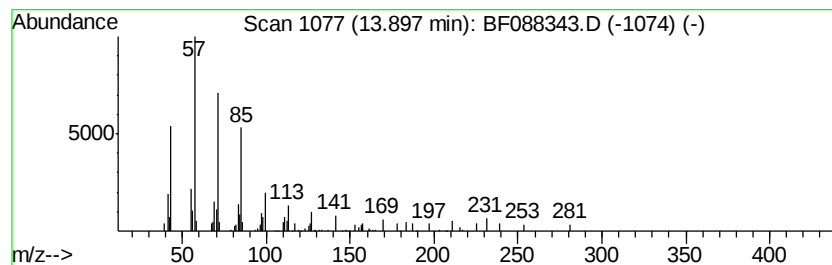
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Peak Number 8 Triacontane Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.90	2.48 ng	65108	Chrysene-d12	13.71

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Triacontane	422	C30H62	000638-68-6	91
2		Heptacosane	380	C27H56	000593-49-7	90
3		Hexacosane	366	C26H54	000630-01-3	90
4		Octacosane	394	C28H58	000630-02-4	90
5		Pentacosane	352	C25H52	000629-99-2	90



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ALS Vial : 13 Sample Multiplier: 1

Instrument :
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ClientSampleId :
Q8-B3(0-8)C

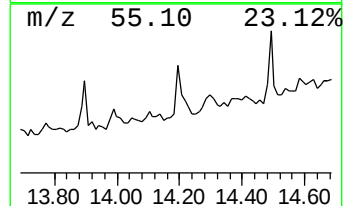
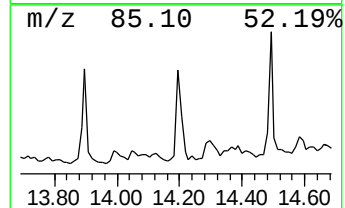
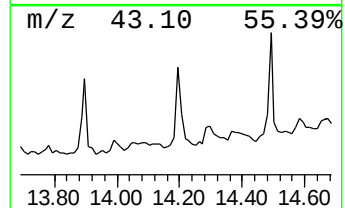
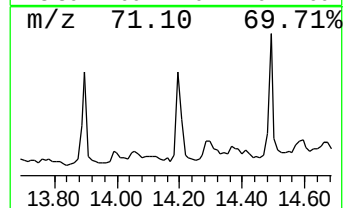
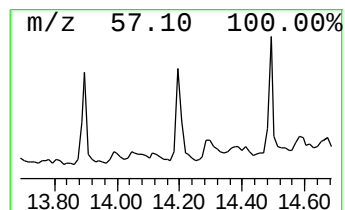
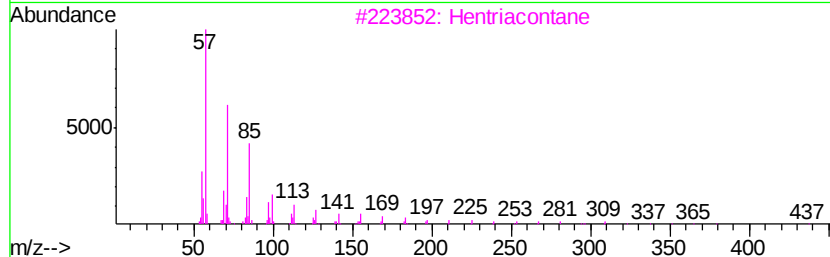
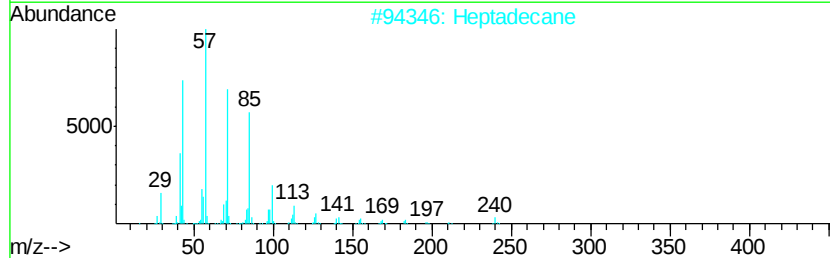
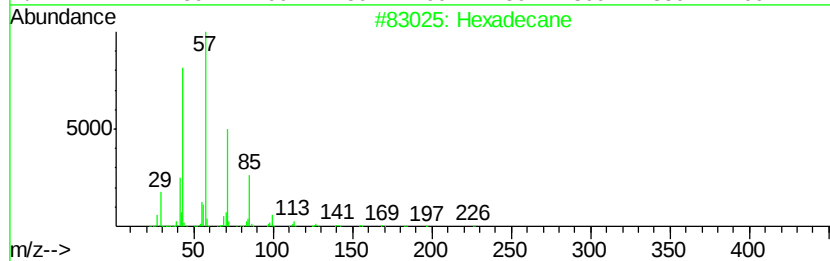
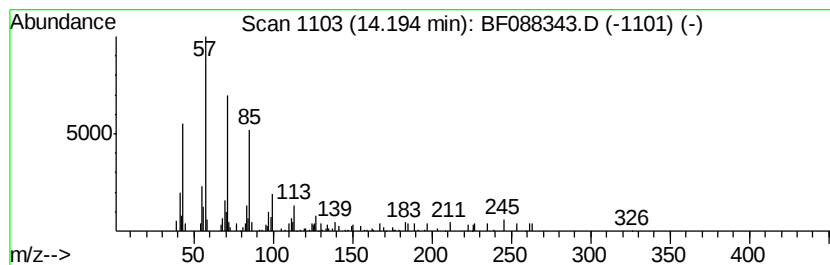
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Hexadecane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.19	3.87 ng	101621	Chrysene-d12	13.71

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane		226	C16H34	000544-76-3	95
2	Heptadecane		240	C17H36	000629-78-7	83
3	Hentriacontane		437	C31H64	000630-04-6	83
4	Heneicosane		296	C21H44	000629-94-7	83
5	Hexacosane		366	C26H54	000630-01-3	83



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF062416\
Data File : BF088343.D
Acq On : 24 Jun 2016 20:49
Operator : UM/SJ
Sample : H3602-15 10X
Misc :
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Instrument :
BNA_F
ClientSampleId :
Q8-B3(0-8)C

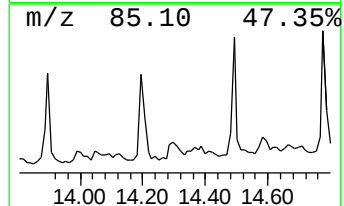
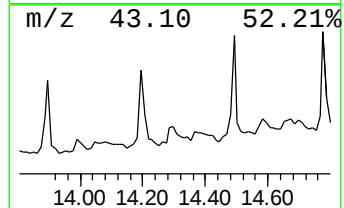
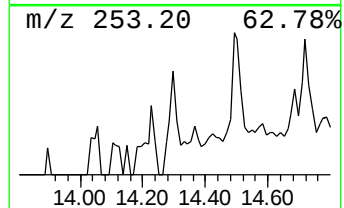
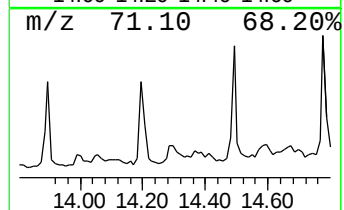
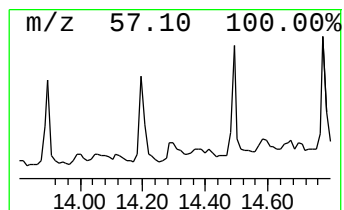
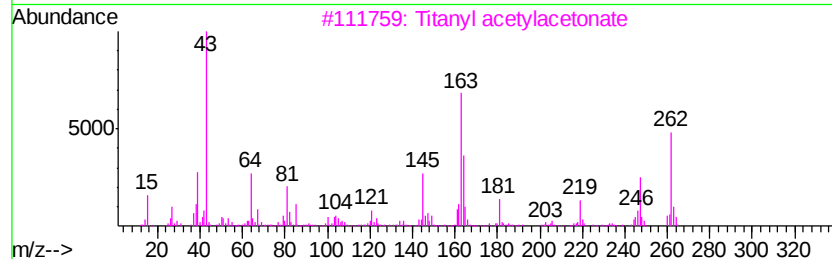
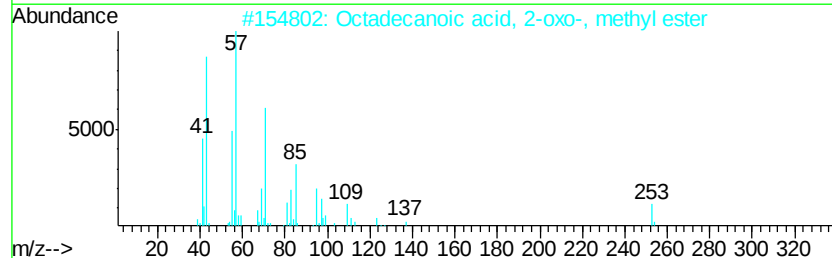
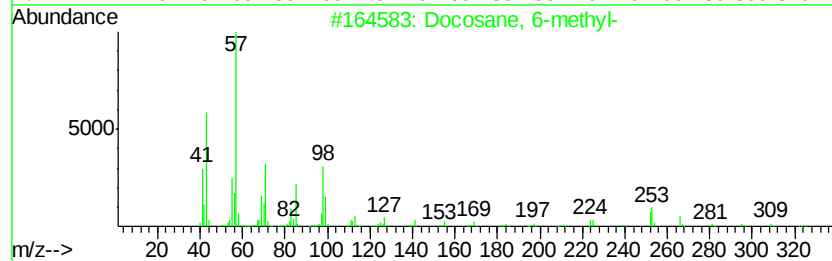
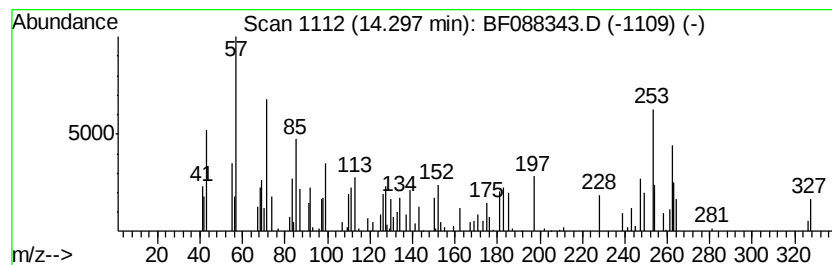
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 unknown14.30 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.30	2.95 ng	77492	Chrysene-d12	13.71

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Docosane, 6-methyl-	324	C23H48	055124-81-7	16
2			Octadecanoic acid, 2-oxo-, methyl...	312	C19H36O3	002380-18-9	14
3			Titanyl acetylacetonate	262	C10H14O5Ti	014024-64-7	12
4			Succinic acid, 2-isopropoxyphenyl...	404	C25H24O5	1000357-97-7	10
5			cis-2,3,6-Trimethoxy-.beta.-meth...	253	C12H15NO5	005556-80-9	10



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF062416\
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Sample : H3602-15 10X
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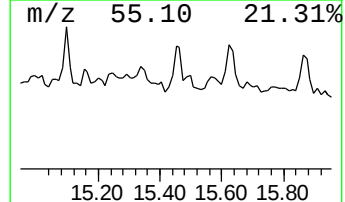
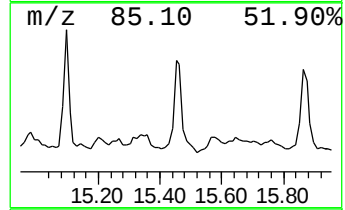
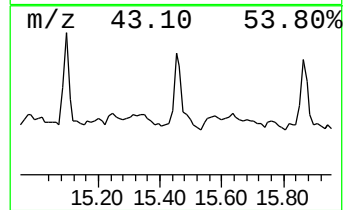
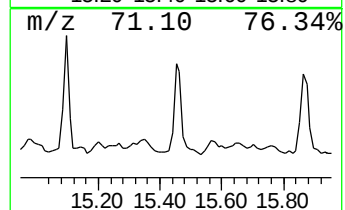
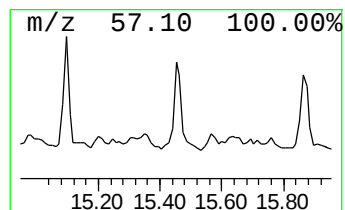
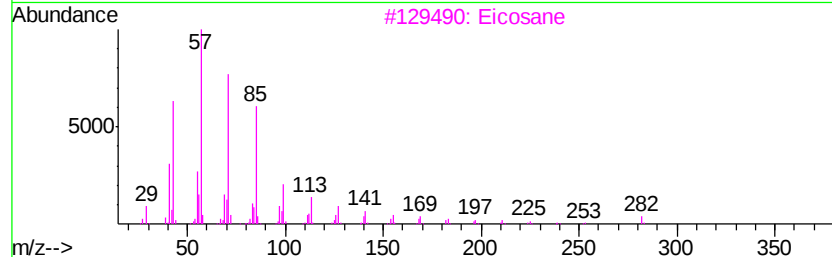
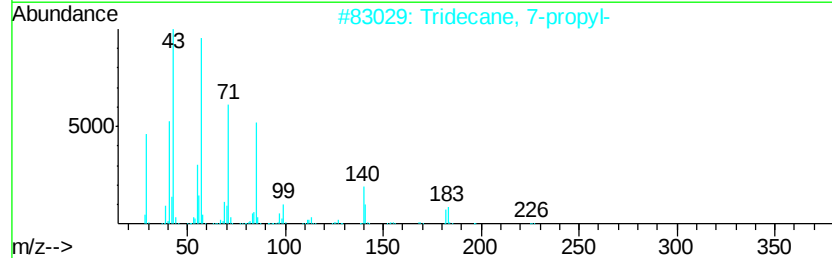
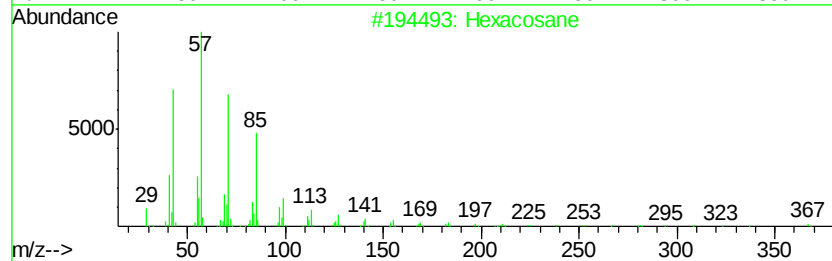
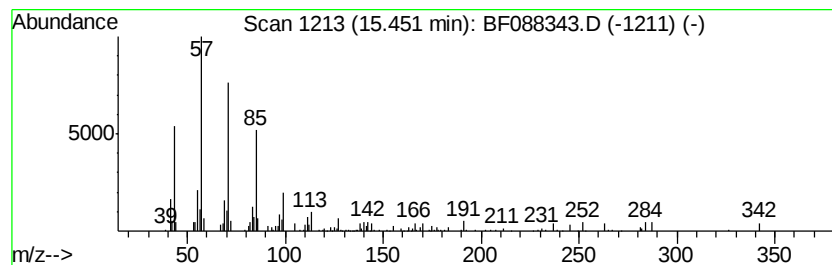
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Hexacosane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.45	2.54 ng	68787	Perylene-d12	15.07
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Hexacosane	366 C26H54	000630-01-3	76
2	Tridecane, 7-propyl-	226 C16H34	055045-09-5	76
3	Eicosane	282 C20H42	000112-95-8	76
4	Nonadecane, 9-methyl-	282 C20H42	013287-24-6	74
5	Hexadecane	226 C16H34	000544-76-3	74



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF062416\
Data File : BF088343.D
Acq On : 24 Jun 2016 20:49
Operator : UM/SJ
Sample : H3602-15 10X
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
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ClientSampleId :
Q8-B3(0-8)C

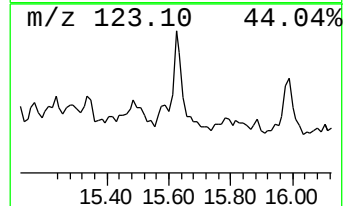
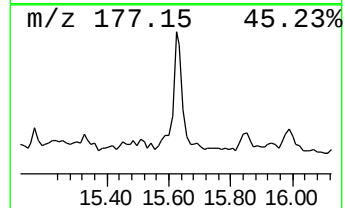
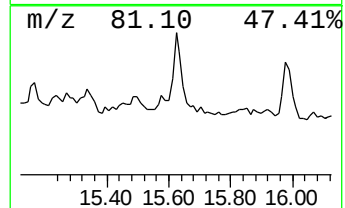
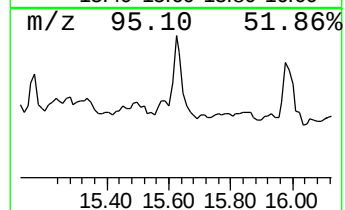
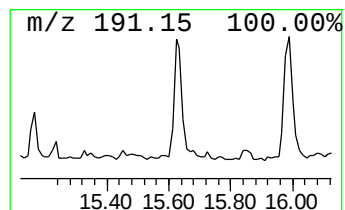
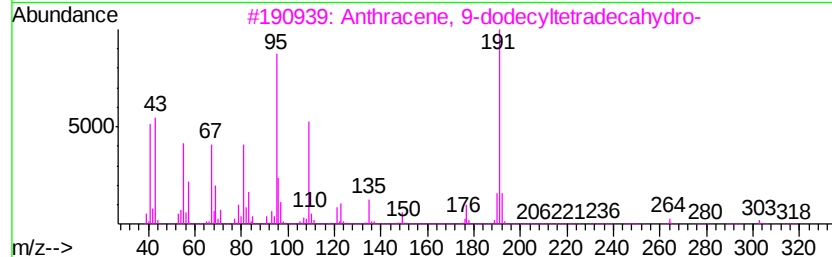
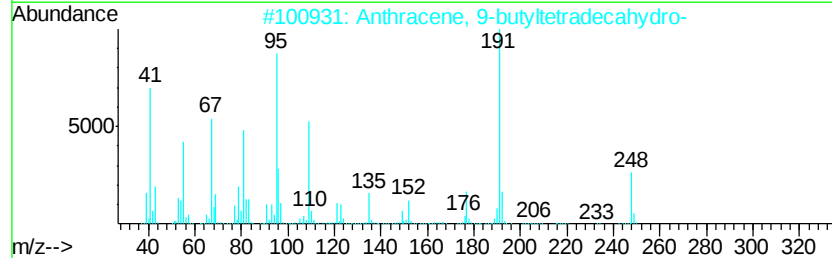
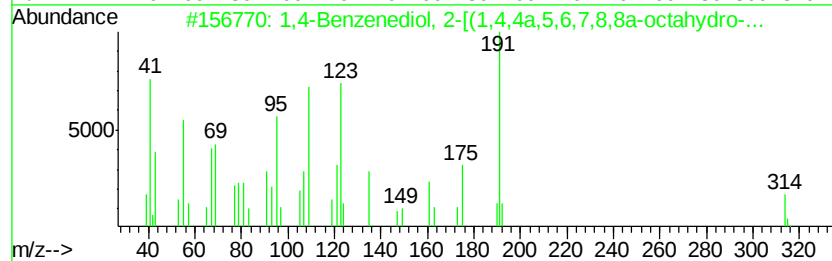
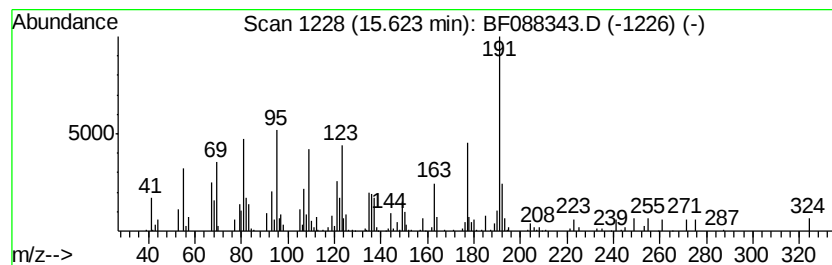
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 unknown15.62 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.62	5.07 ng	137111	Perylene-d12	15.07

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,4-Benzenediol, 2-[(1,4,4a,5,6,...	314	C21H30O2	039707-55-6	46
2			Anthracene, 9-butyltetradeca-	248	C18H32	055133-89-6	38
3			Anthracene, 9-dodecyltetradeca-	360	C26H48	055401-75-7	37
4			2H-1,2,3-Triazole-4-carboxaldehy...	191	C9H6FN3O	051306-43-5	32
5			.beta.-iso-Methyl ionone	206	C14H22O	1000285-40-2	27



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF062416\
Data File : BF088343.D
Acq On : 24 Jun 2016 20:49
Operator : UM/SJ
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Misc :
ALS Vial : 13 Sample Multiplier: 1

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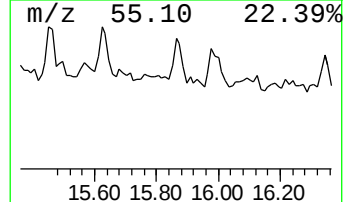
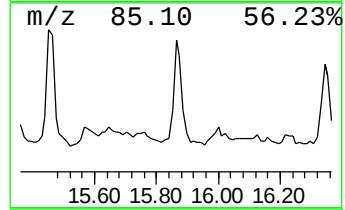
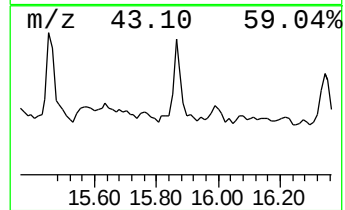
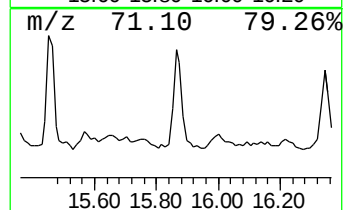
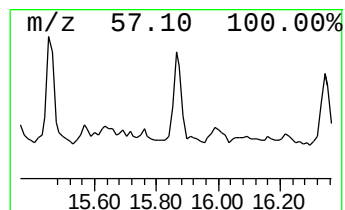
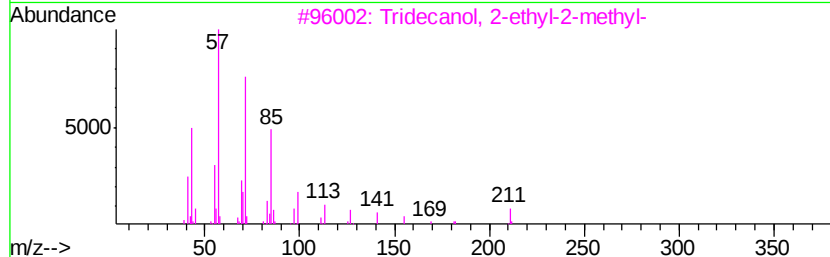
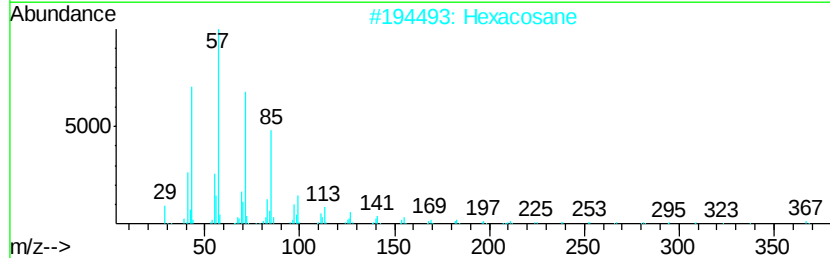
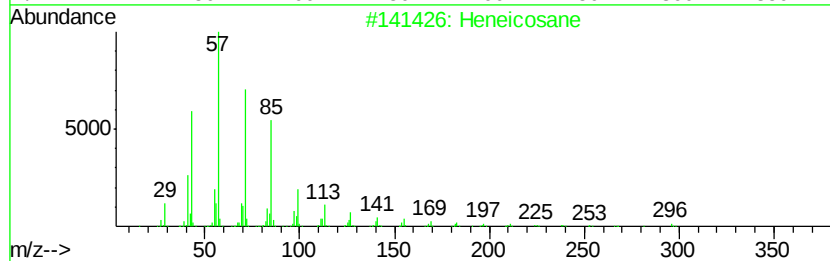
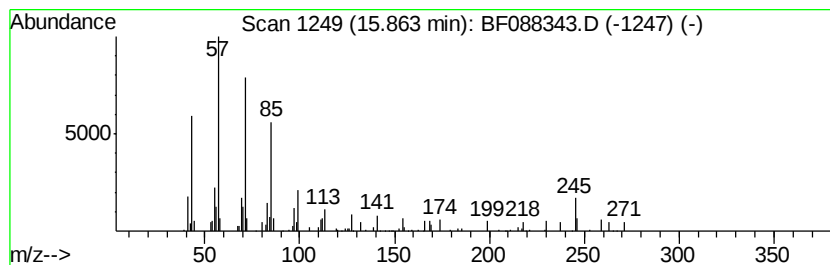
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Heneicosane Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.86	2.43 ng	65821	Perylene-d12	15.07

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	87
2		Hexacosane	366	C26H54	000630-01-3	87
3		Tridecanol, 2-ethyl-2-methyl-	242	C16H34O	1000115-66-1	80
4		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	80
5		Eicosane, 10-methyl-	296	C21H44	054833-23-7	80



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF062416\
Data File : BF088343.D
Acq On : 24 Jun 2016 20:49
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Sample : H3602-15 10X
Misc :
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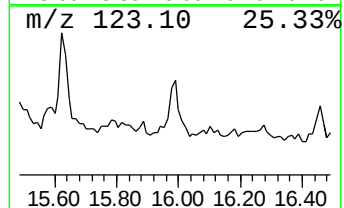
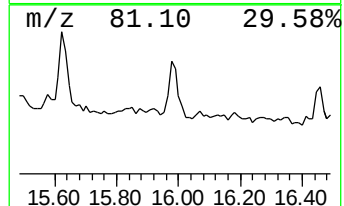
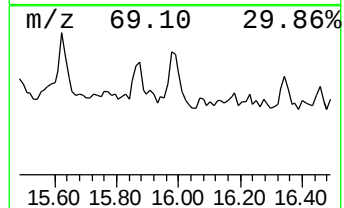
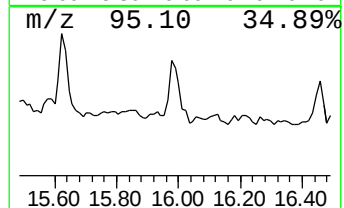
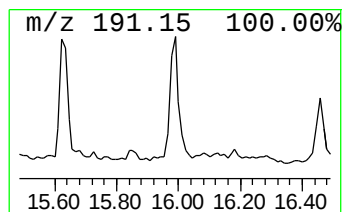
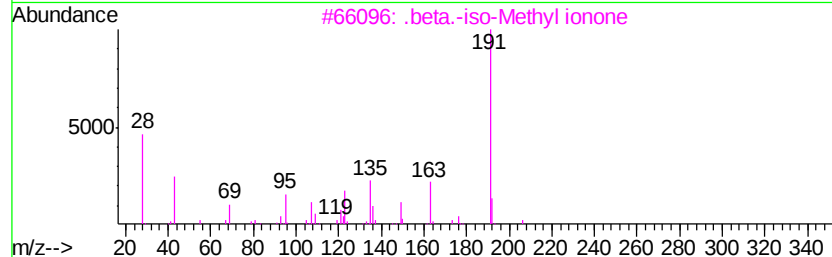
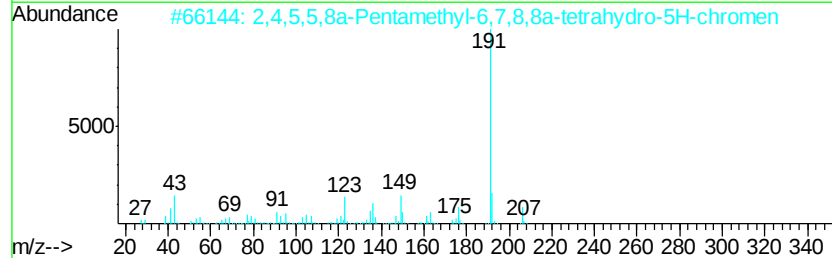
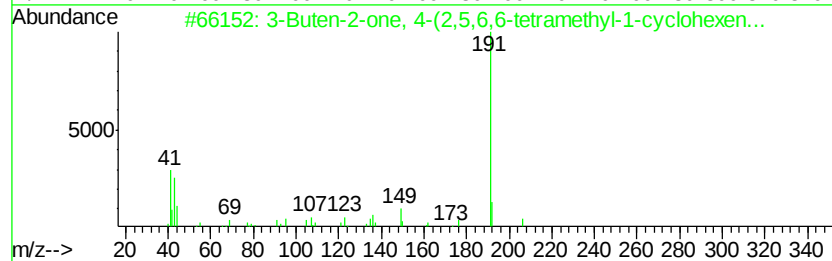
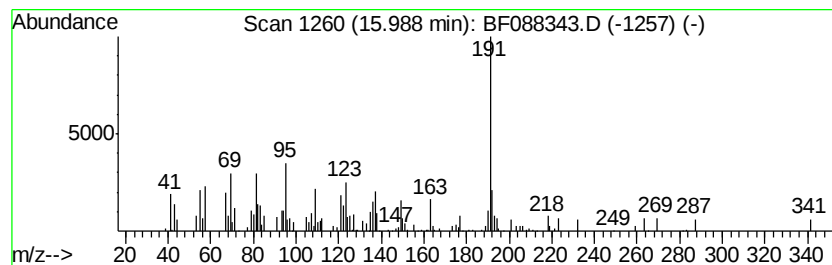
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 unknown15.99 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.	
15.99	4.05 ng	109538	Perylene-d12	15.07	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Buten-2-one, 4-(2,5,6,6-tetram...	206	C14H22O	000079-70-9	49
2	2,4,5,5,8a-Pentamethyl-6,7,8,8a-...	206	C14H22O	1000195-40-9	47
3	.beta.-iso-Methyl ionone	206	C14H22O	1000285-40-2	46
4	Anthracene, 9-(3-butenyl)-	232	C18H16	097451-55-3	38
5	1-(3-Nitro-phenyl)-pyrrolidine	192	C10H12N2O2	1000300-34-7	38



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF062416\
Data File : BF088343.D
Acq On : 24 Jun 2016 20:49
Operator : UM/SJ
Sample : H3602-15 10X
Misc :
ALS Vial : 13 Sample Multiplier: 1

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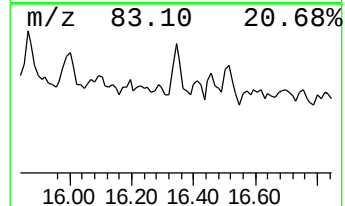
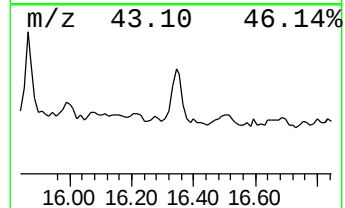
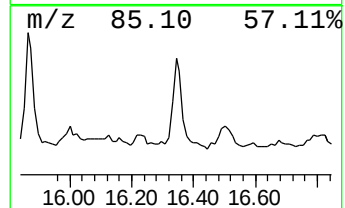
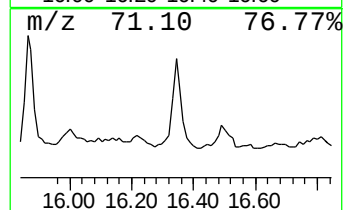
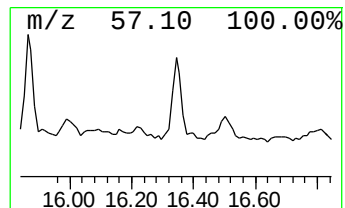
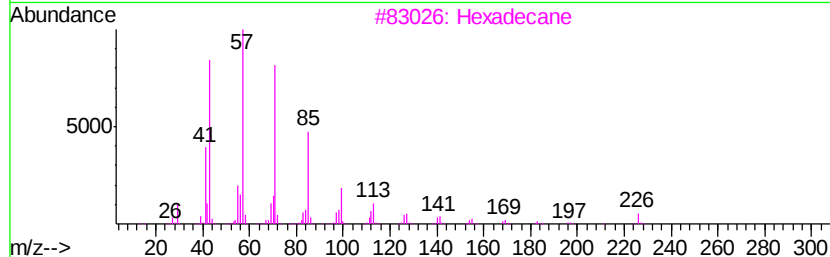
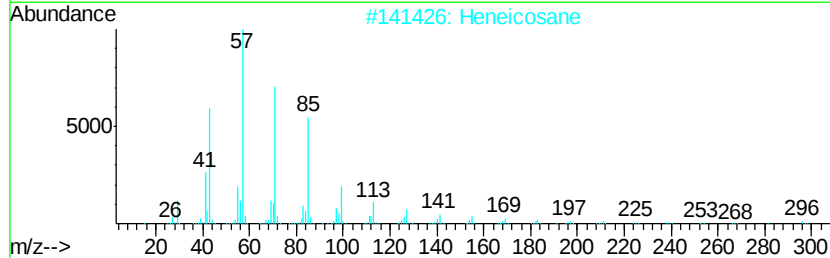
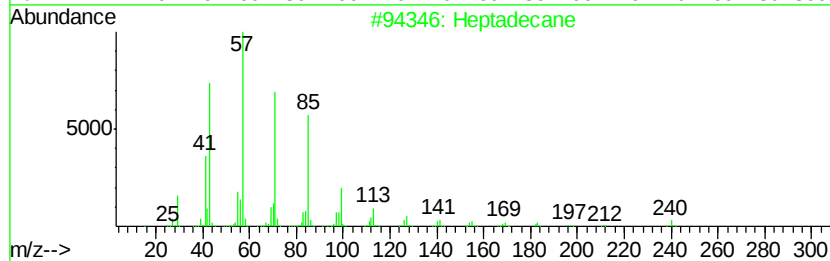
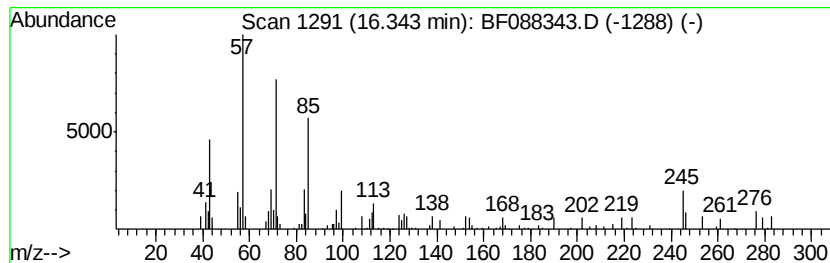
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 Heptadecane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.34	2.63 ng	71162	Perylene-d12	15.07

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptadecane	240	C17H36	000629-78-7	64
2		Heneicosane	296	C21H44	000629-94-7	64
3		Hexadecane	226	C16H34	000544-76-3	64
4		Octadecane, 1-iodo-	380	C18H37I	000629-93-6	64
5		Octadecane, 2-methyl-	268	C19H40	001560-88-9	59



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF062416\
Data File : BF088343.D
Acq On : 24 Jun 2016 20:49
Operator : UM/SJ
Sample : H3602-15 10X
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleID :
Q8-B3(0-8)C

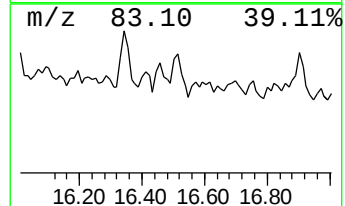
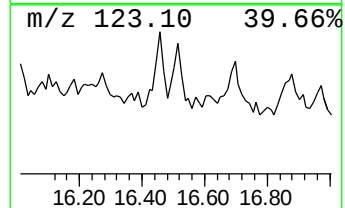
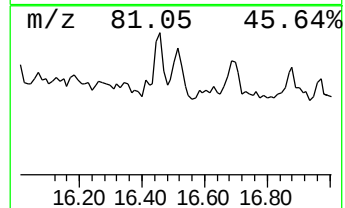
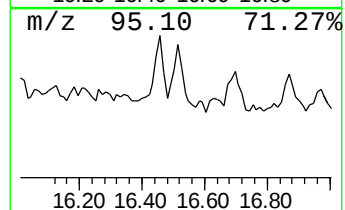
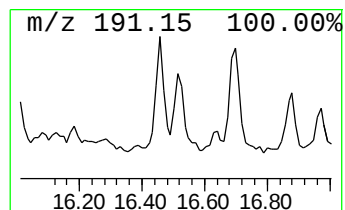
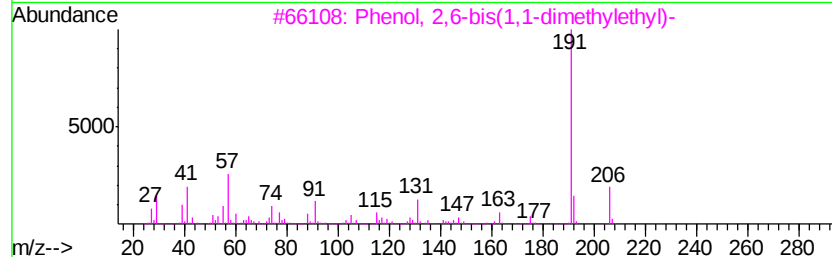
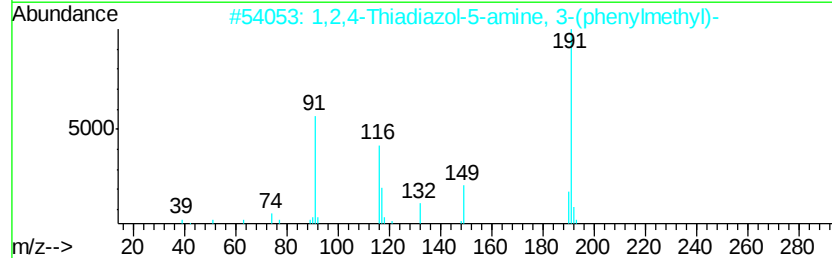
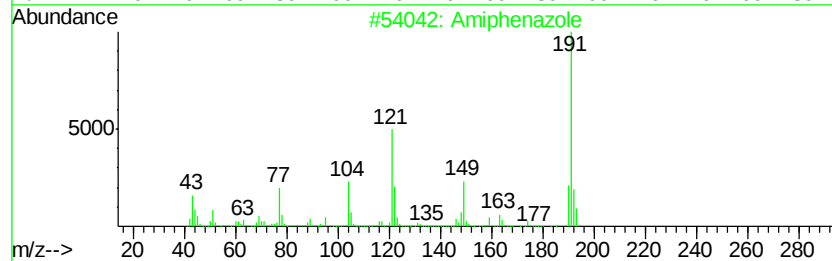
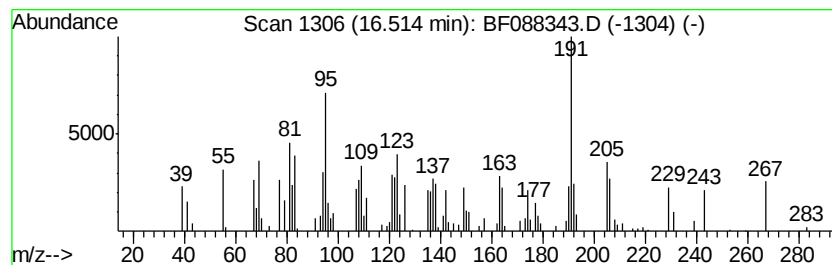
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060716.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 unknown16.51 Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.51	2.38 ng	64314	Perylene-d12	15.07

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Amiphenazole	191	C9H9N3S	000490-55-1	38
2		1,2,4-Thiadiazol-5-amine, 3-(phe...	191	C9H9N3S	017467-27-5	25
3		Phenol, 2,6-bis(1,1-dimethylethyl)-	206	C14H22O	000128-39-2	22
4		1-(3,6,6-Trimethyl-1,6,7,7a-tetr...	206	C13H18O2	1000194-97-2	22
5		2,5,5,8a-Tetramethyl-4-methylene...	206	C14H22O	093175-76-9	22



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF062416\
Data File : BF088343.D
Acq On : 24 Jun 2016 20:49
Operator : UM/SJ
Sample : H3602-15 10X
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
Q8-B3(0-8)C

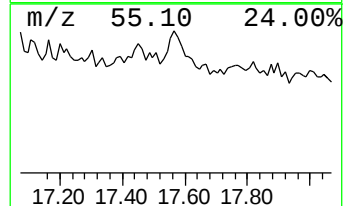
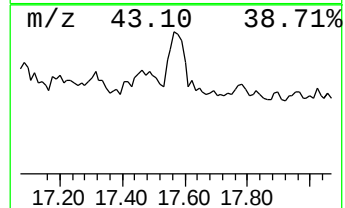
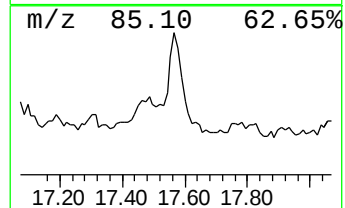
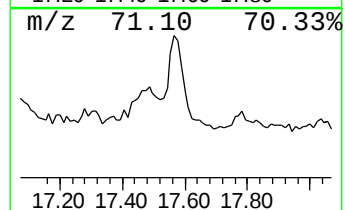
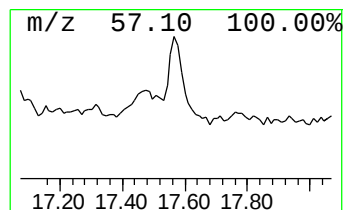
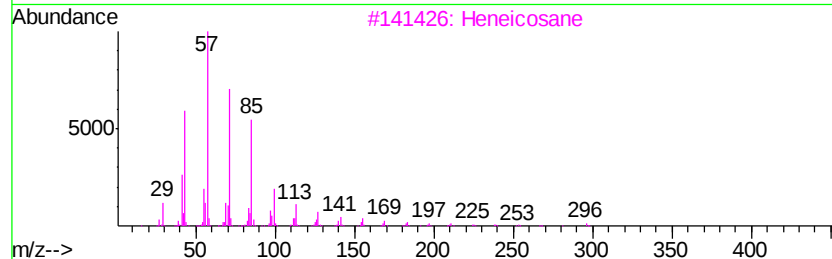
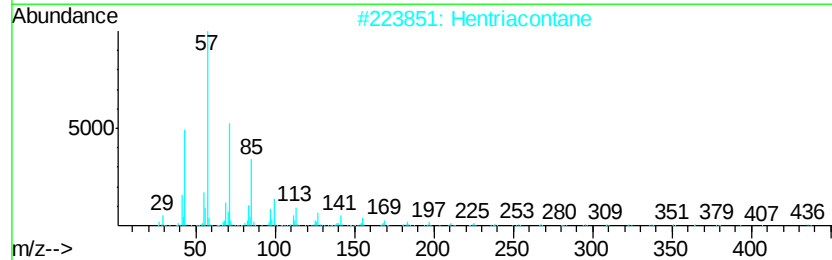
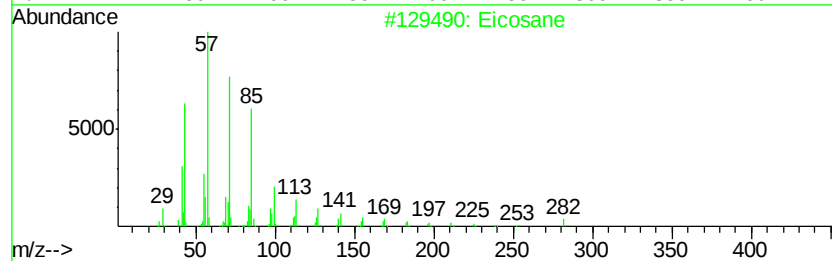
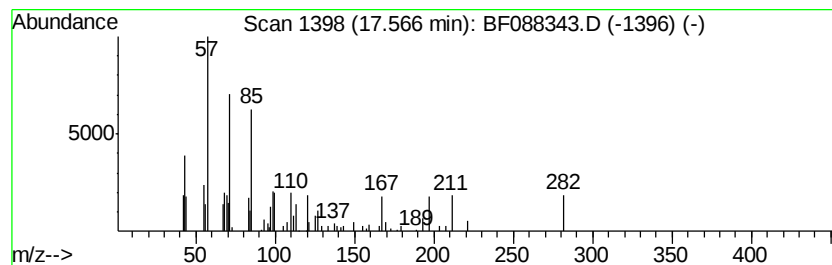
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060716.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 Eicosane Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.57	2.22 ng	59971	Perylene-d12	15.07

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	64
2		Hentriacontane	437	C31H64	000630-04-6	49
3		Heneicosane	296	C21H44	000629-94-7	49
4		Hexacosane	366	C26H54	000630-01-3	49
5		Octacosane	394	C28H58	000630-02-4	49



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF062416\
Data File : BF088343.D
Acq On : 24 Jun 2016 20:49
Operator : UM/SJ
Sample : H3602-15 10X
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
Q8-B3(0-8)C

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060716.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown6.35	6.35	10.1 ng		214464	1	6.58	425723	20.0
Hexadecanoic acid...	11.50	25.2 ng		710824	4	11.09	564933	20.0
9-Octadecenoic ac...	12.21	115.1 ng		3251160	4	11.09	564933	20.0
Methyl stearate	12.29	9.8 ng		277333	4	11.09	564933	20.0
Triacontane	13.90	2.5 ng		65108	5	13.71	525352	20.0
Hexadecane	14.19	3.9 ng		101621	5	13.71	525352	20.0
unknown14.30	14.30	3.0 ng		77492	5	13.71	525352	20.0
Hexacosane	15.45	2.5 ng		68787	6	15.07	541024	20.0
unknown15.62	15.62	5.1 ng		137111	6	15.07	541024	20.0
Heneicosane	15.86	2.4 ng		65821	6	15.07	541024	20.0
unknown15.99	15.99	4.0 ng		109538	6	15.07	541024	20.0
Heptadecane	16.34	2.6 ng		71162	6	15.07	541024	20.0
unknown16.51	16.51	2.4 ng		64314	6	15.07	541024	20.0
Eicosane	17.57	2.2 ng		59971	6	15.07	541024	20.0