

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF011116\
 Data File : BF084379.D
 Acq On : 11 Jan 2016 14:58
 Operator : SJ/UM
 Sample : H1043-18
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-4COMP

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-BF123115.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	5.047	256	259	267	rVB	1385053	2184009	25.43%	3.163%
2	5.470	293	296	299	rBV	6135748	7202862	83.86%	10.431%
3	6.510	383	387	389	rBV	6771663	8588906	100.00%	12.438%
4	6.624	395	397	400	rBV	5652428	7767760	90.44%	11.249%
5	6.853	415	417	419	rBV	1966682	1587024	18.48%	2.298%
6	7.013	428	431	433	rBV	6401246	5744752	66.89%	8.319%
7	7.425	463	467	469	rBV	4984340	5224759	60.83%	7.566%
8	8.145	527	530	532	rBV	1938713	2053304	23.91%	2.973%
9	9.219	621	624	627	rBV	7395669	6946904	80.88%	10.060%
10	9.619	656	659	660	rBV	1346490	1545025	17.99%	2.237%
11	9.825	676	677	680	rBV	171669	213950	2.49%	0.310%
12	9.893	681	683	686	rVB	2362929	2080337	24.22%	3.013%
13	10.328	720	721	723	rVB	456400	363922	4.24%	0.527%
14	10.545	737	740	744	rBV	116488	231934	2.70%	0.336%
15	10.682	749	752	755	rVV2	3785699	4578467	53.31%	6.630%
16	10.796	759	762	770	rVV2	457234	815253	9.49%	1.181%
17	10.991	777	779	784	rBV	50407	116055	1.35%	0.168%
18	11.253	800	802	803	rBV	218694	208142	2.42%	0.301%
19	11.379	811	813	815	rBV	2165035	1883228	21.93%	2.727%
20	12.785	934	936	938	rBV	259977	314888	3.67%	0.456%
21	12.854	938	942	949	rVB2	169363	304784	3.55%	0.441%
22	12.968	949	952	954	rBV	5855507	5600996	65.21%	8.111%
23	13.494	996	998	1000	rBV	251894	221604	2.58%	0.321%
24	13.871	1028	1031	1037	rBV	162495	267776	3.12%	0.388%
25	14.008	1041	1043	1046	rBV	1007145	1273879	14.83%	1.845%
26	15.437	1165	1168	1171	rBV	752229	1136756	13.24%	1.646%
27	16.374	1247	1250	1259	rVB2	49783	109158	1.27%	0.158%
28	16.945	1296	1300	1304	rBV	47694	108397	1.26%	0.157%
29	17.608	1354	1358	1362	rVB	53198	128958	1.50%	0.187%
30	18.408	1423	1428	1433	rBV2	41135	122727	1.43%	0.178%
31	19.346	1506	1510	1515	rVB3	37127	127618	1.49%	0.185%

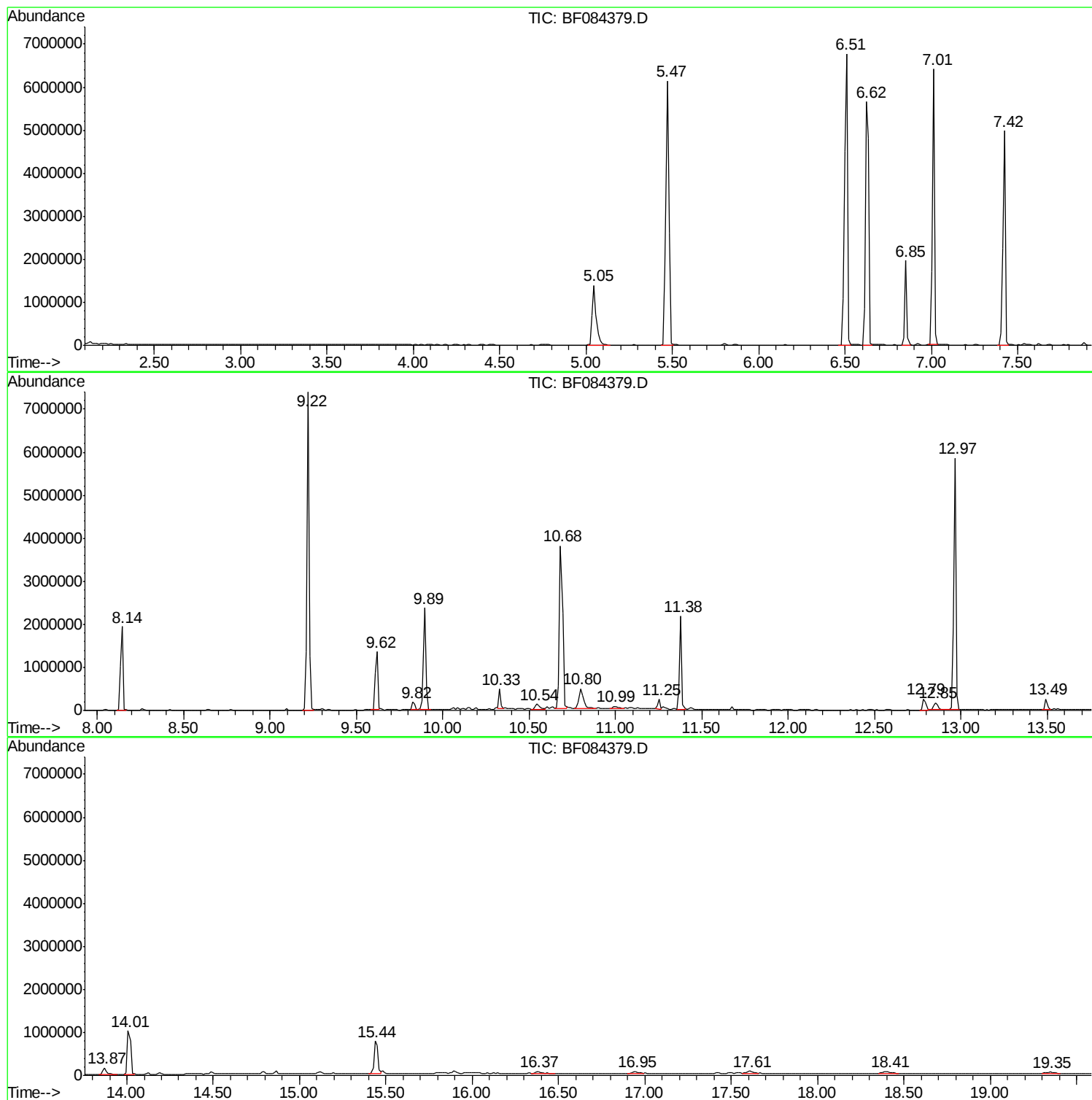
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Instrument :
BNA_F
ClientSampleId :
SB-4COMP

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF123115.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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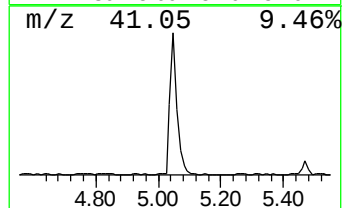
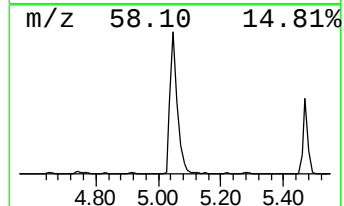
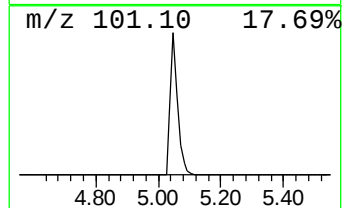
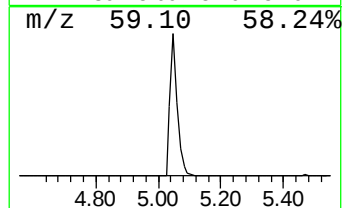
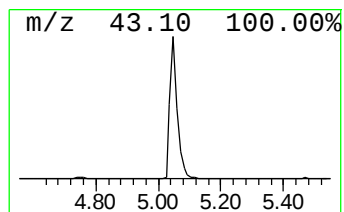
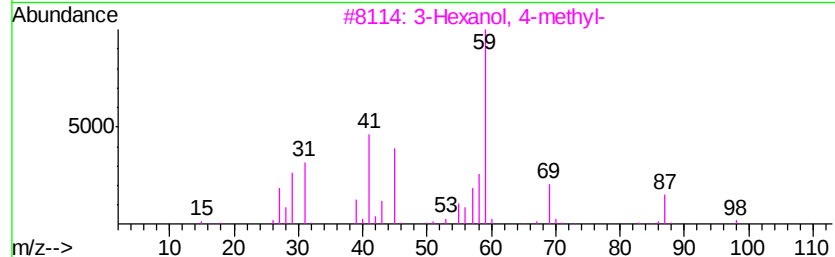
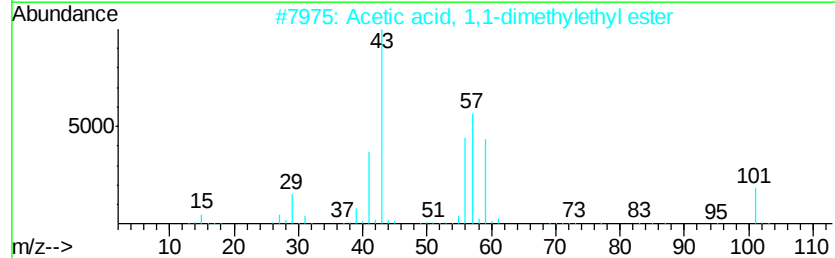
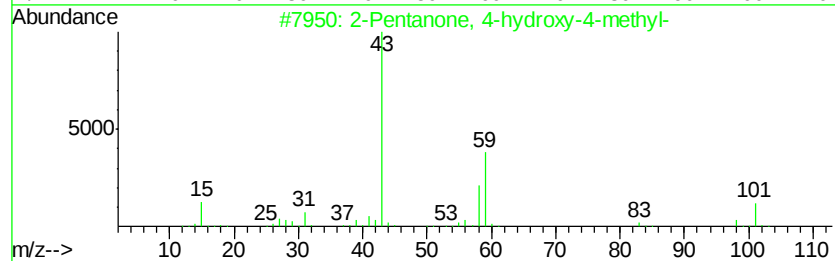
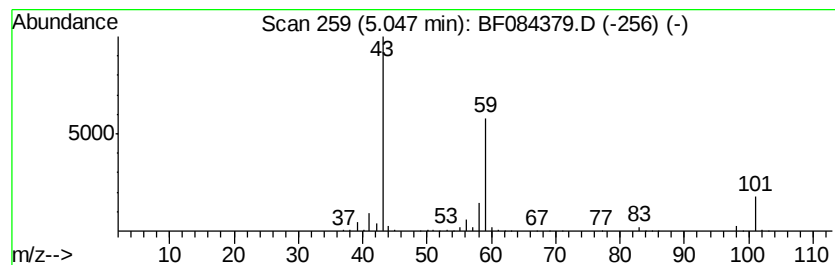
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.05	27.52 ng	2184010	1,4-Dichlorobenzene-d4	6.85

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	53
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			Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	33
5			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	32



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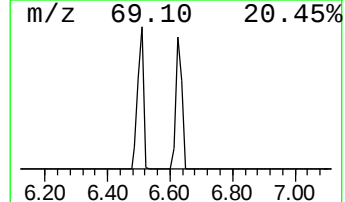
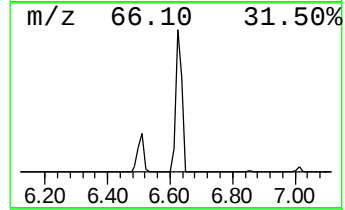
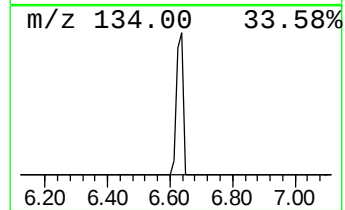
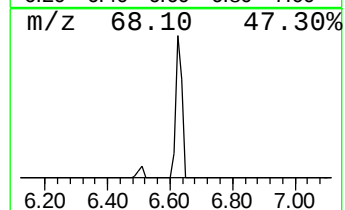
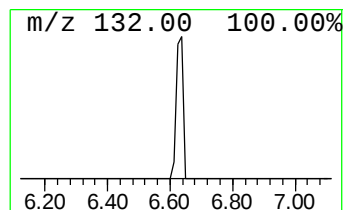
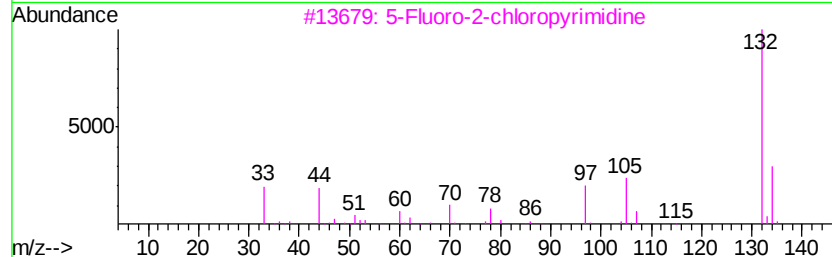
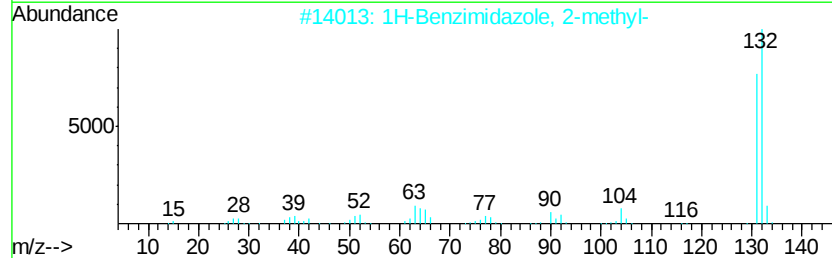
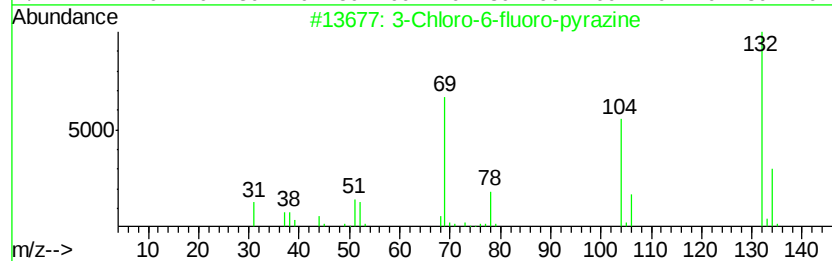
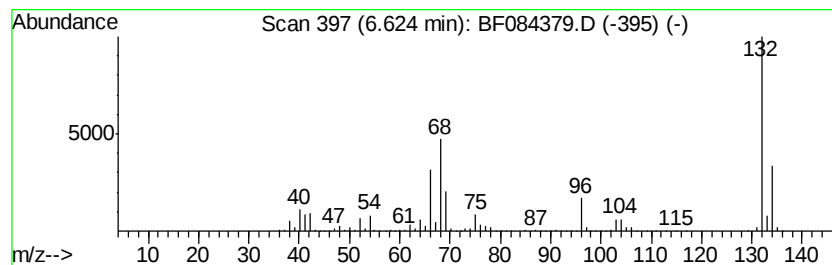
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Peak Number 2 unknown6.62 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.62	97.89 ng	7767760	1,4-Dichlorobenzene-d4	6.85

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Chloro-6-fluoro-pyrazine			132	C4H2ClFN2	1000146-10-7	27
2	1H-Benzimidazole, 2-methyl-			132	C8H8N2	000615-15-6	22
3	5-Fluoro-2-chloropyrimidine			132	C4H2ClFN2	062802-42-0	12
4	(5-METHYL-2-PYRIDYL)ACETONITRILE			132	C8H8N2	1000241-93-9	10
5	1,2,3-Trifluorobenzene			132	C6H3F3	001489-53-8	9



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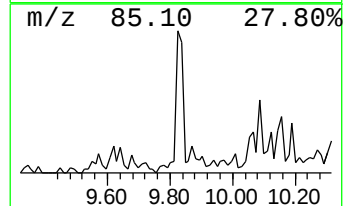
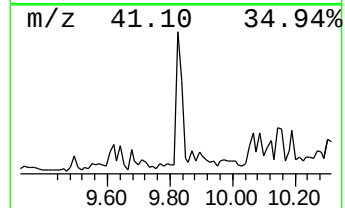
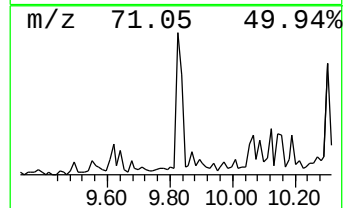
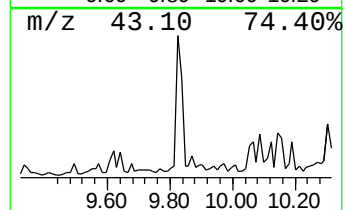
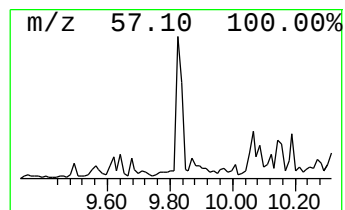
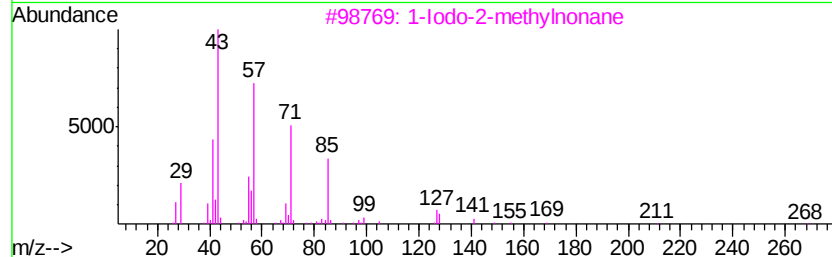
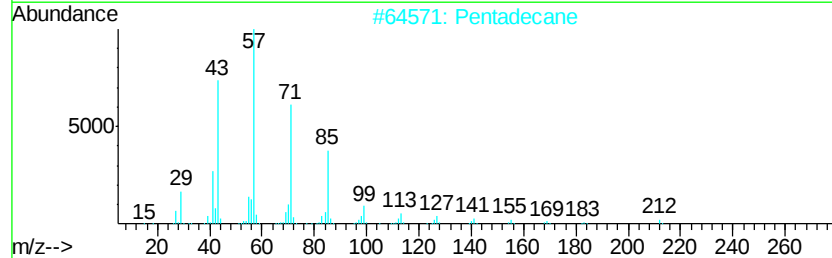
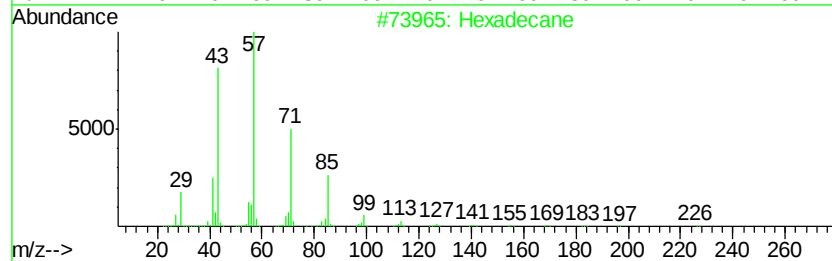
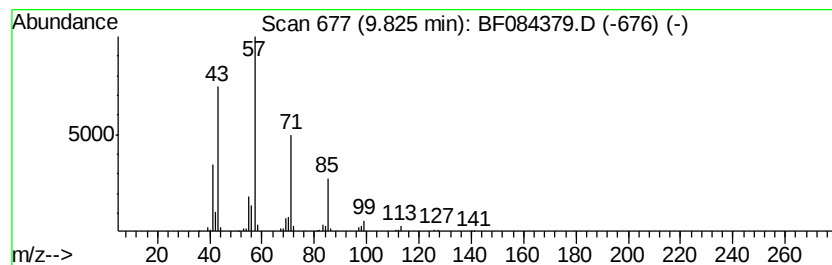
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Hexadecane Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.82	2.06 ng	213950	Acenaphthene-d10	9.89
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Hexadecane	226 C16H34	000544-76-3	91
2	Pentadecane	212 C15H32	000629-62-9	78
3	1-Iodo-2-methylnonane	268 C10H21I	1000101-47-9	59
4	Tetradecane	198 C14H30	000629-59-4	59
5	Decane, 2,5,9-trimethyl-	184 C13H28	062108-22-9	59



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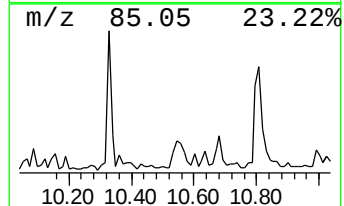
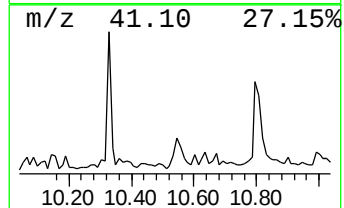
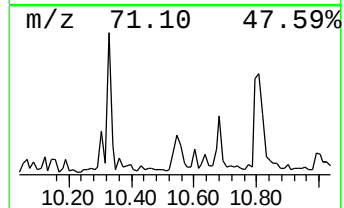
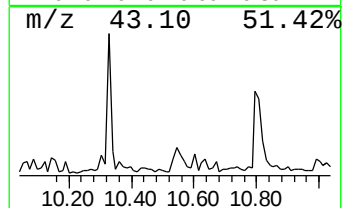
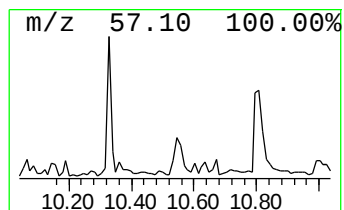
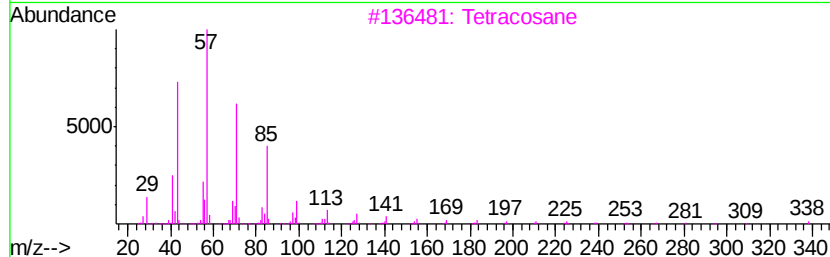
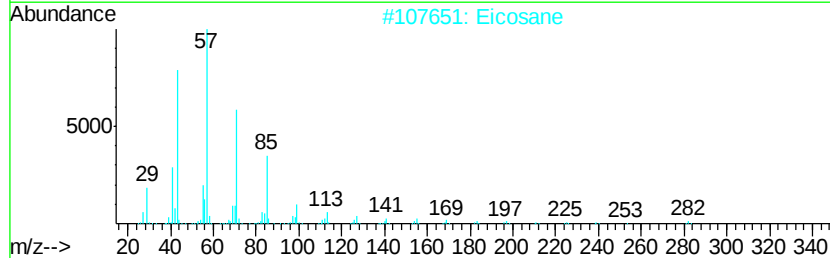
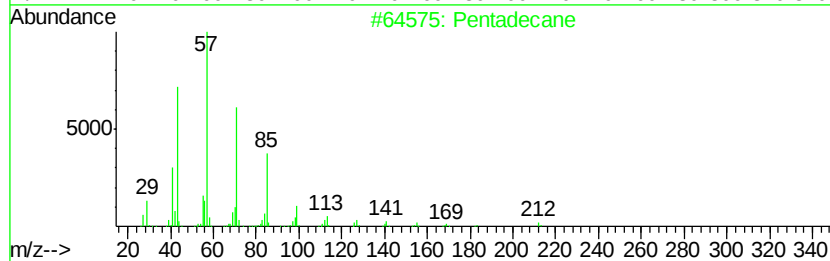
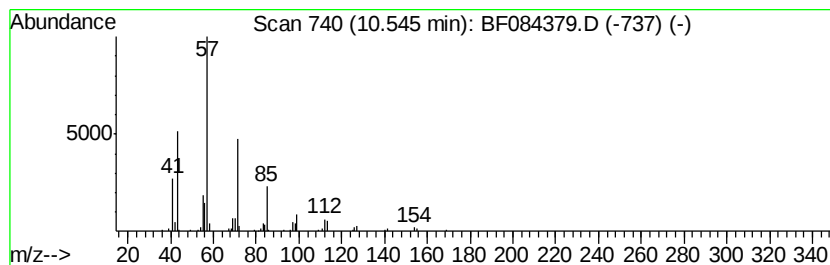
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 6 Pentadecane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.54	2.23 ng	231934	Acenaphthene-d10	9.89
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Pentadecane	212 C15H32	000629-62-9	80
2	Eicosane	282 C20H42	000112-95-8	72
3	Tetracosane	338 C24H50	000646-31-1	72
4	Heptacosane	380 C27H56	000593-49-7	68
5	Tridecane, 3-methyl-	198 C14H30	006418-41-3	64



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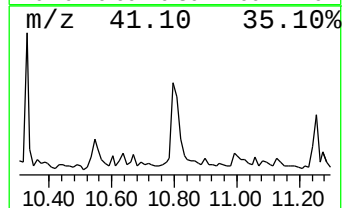
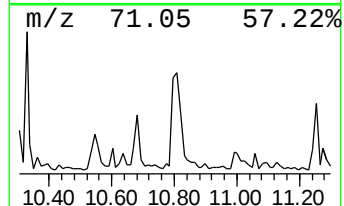
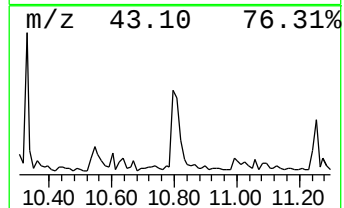
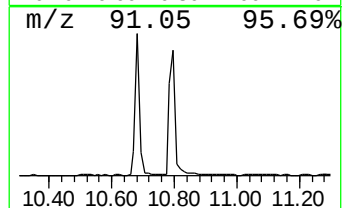
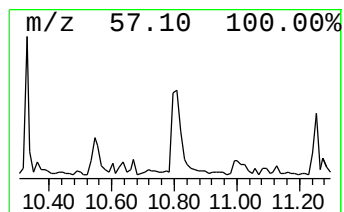
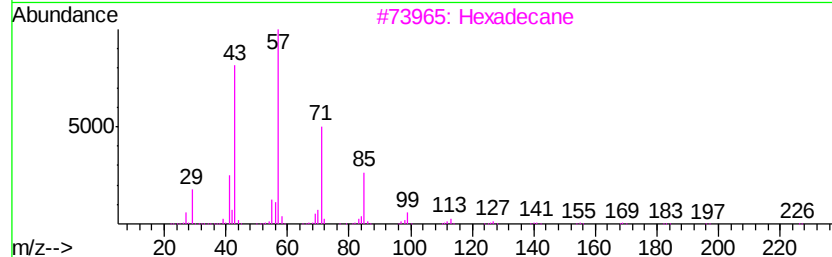
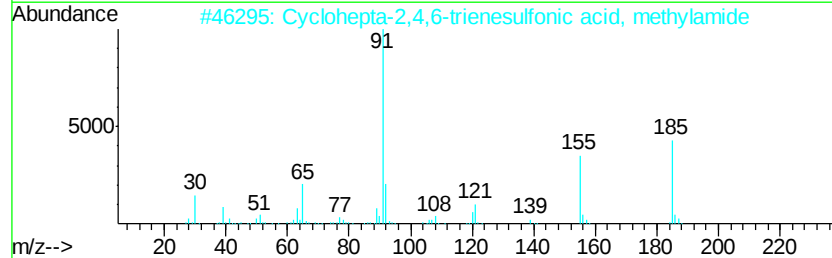
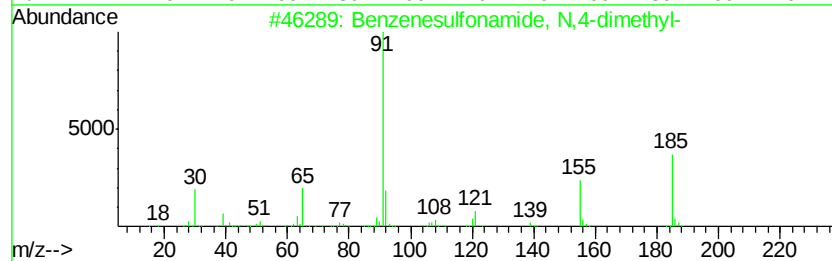
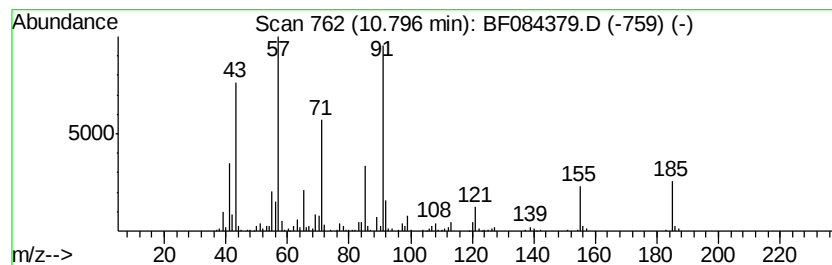
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TIC Integration Parameters: LSCINT.P

Peak Number 7 Benzenesulfonamide, N,4-dim... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.80	8.66 ng	815253	Phenanthrene-d10	11.38	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzenesulfonamide, N,4-dimethyl-	185	C8H11NO2S	000640-61-9	95
2	Cyclohepta-2,4,6-trienesulfonic ...	185	C8H11NO2S	1000187-28-7	80
3	Hexadecane	226	C16H34	000544-76-3	38
4	Tridecane, 6-propyl-	226	C16H34	055045-10-8	35
5	4-Toluenesulfonylmethylisocyanide	195	C9H9NO2S	036635-61-7	22



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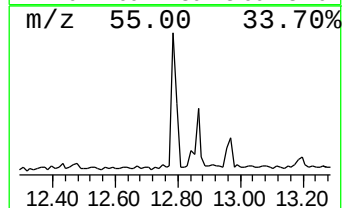
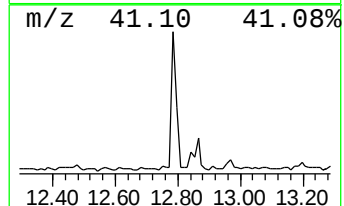
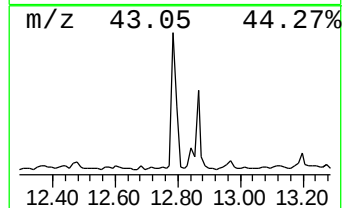
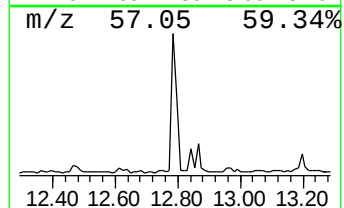
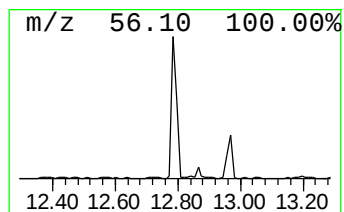
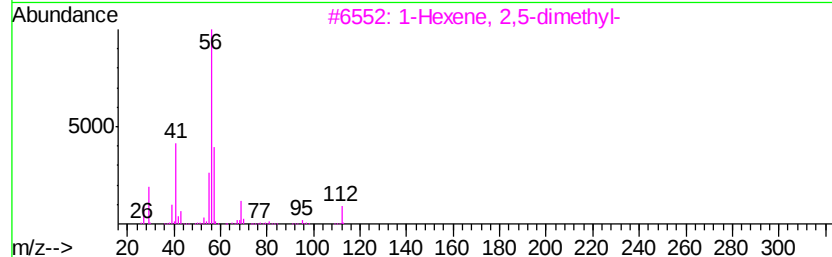
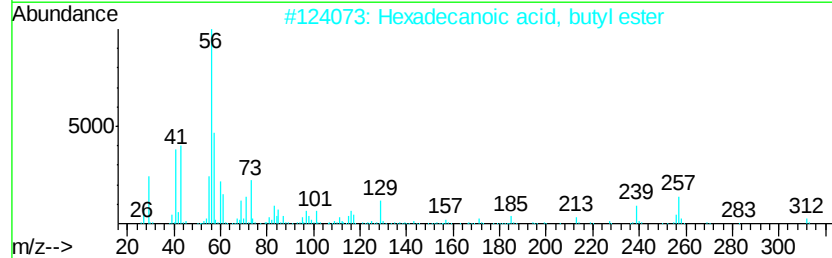
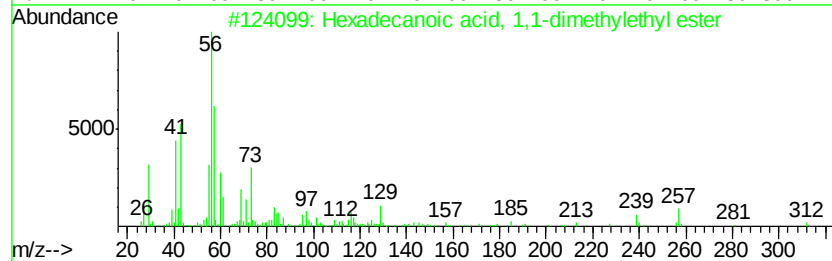
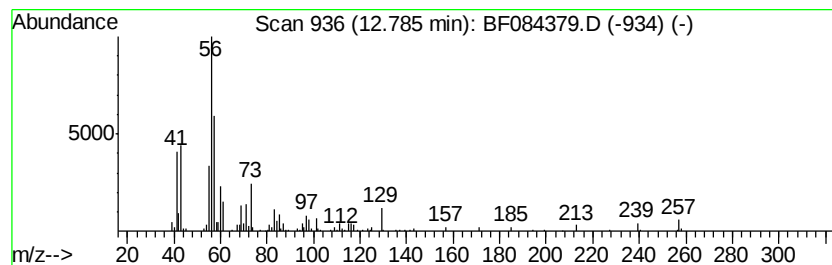
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Hexadecanoic acid, 1,1-dime... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.79	4.94 ng	314888	Chrysene-d12	14.01

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecanoic acid, 1,1-dimethyle...	312	C20H40O2	031158-91-5	90
2		Hexadecanoic acid, butyl ester	312	C20H40O2	000111-06-8	80
3		1-Hexene, 2,5-dimethyl-	112	C8H16	006975-92-4	38
4		Cyclobutane, 1,1-dimethyl-2-octyl-	196	C14H28	062338-30-1	38
5		1-Pentene, 2,4-dimethyl-	98	C7H14	002213-32-3	38



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF011116\
Data File : BF084379.D
Acq On : 11 Jan 2016 14:58
Operator : SJ/UM
Sample : H1043-18
Misc :
ALS Vial : 9 Sample Multiplier: 1

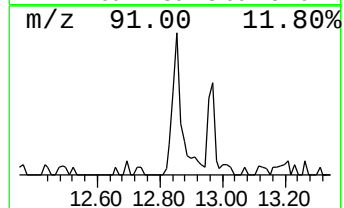
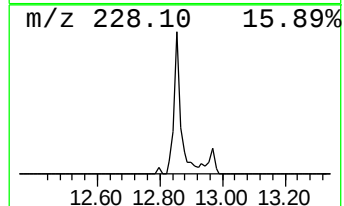
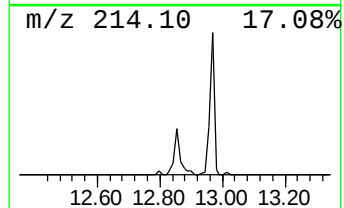
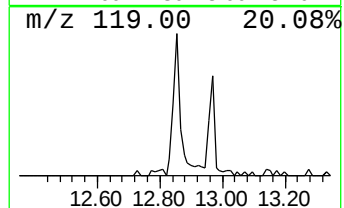
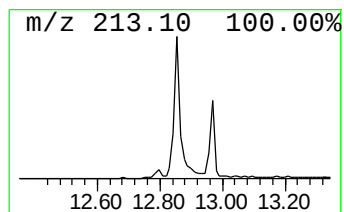
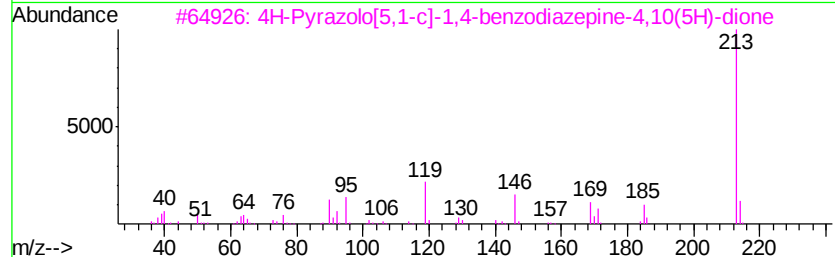
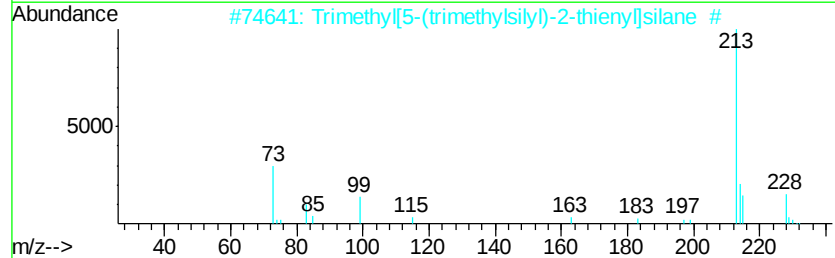
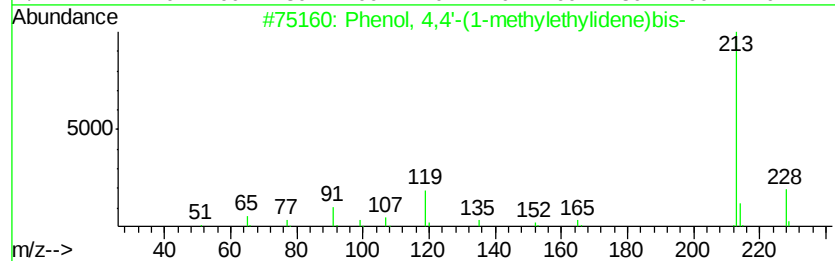
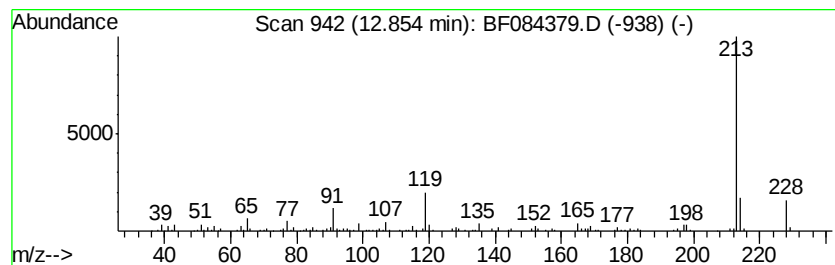
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Phenol, 4,4'-(1-methylethyl... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.85	4.79 ng	304784	Chrysene-d12	14.01	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 4,4'-(1-methylethylidene...	228	C15H16O2	000080-05-7	97
2	Trimethyl[5-(trimethylsilyl)-2-t...	228	C10H20SSi2	017906-71-7	53
3	4H-Pyrazolo[5,1-c]-1,4-benzodiaz...	213	C11H7N3O2	1000277-20-1	45
4	Naphthalene-2,6-dicarboxylic aci...	376	C24H24O4	1000189-60-8	40
5	2H,8H-Benzo[1,2-b:3,4-b']dipyran...	228	C14H12O3	000523-59-1	40



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF011116\
Data File : BF084379.D
Acq On : 11 Jan 2016 14:58
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Instrument :
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ClientSampleId :
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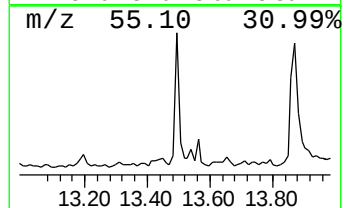
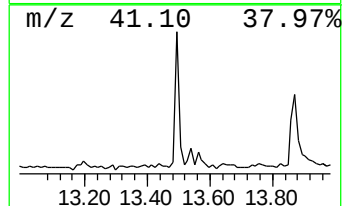
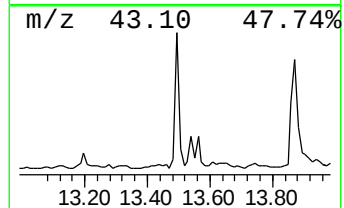
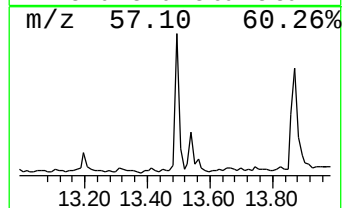
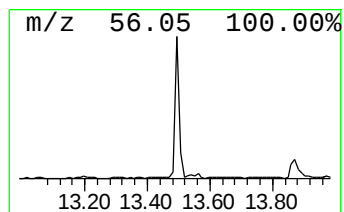
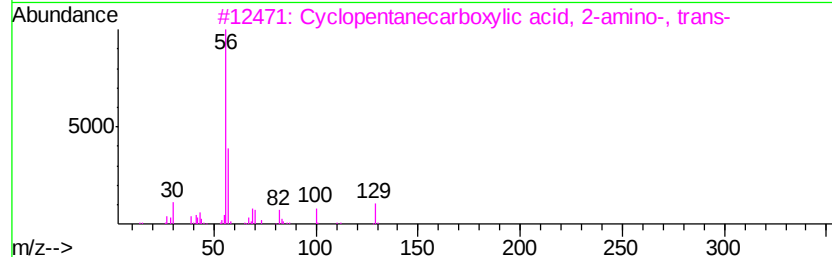
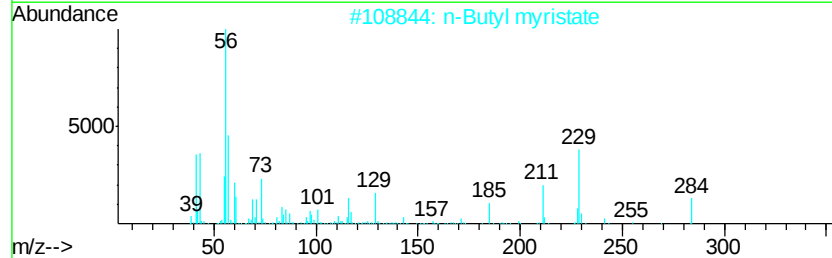
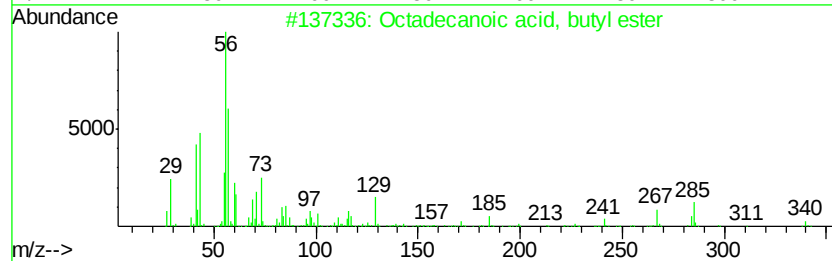
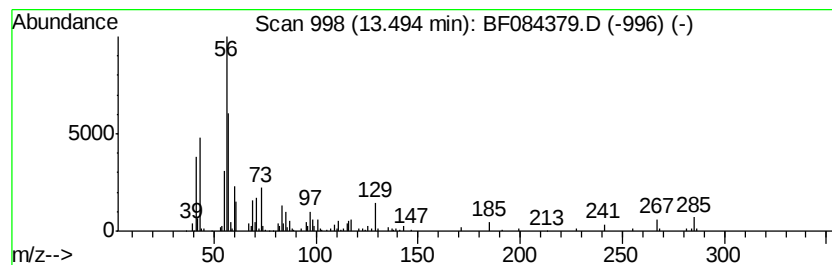
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Octadecanoic acid, butyl ester Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.49	3.48 ng	221604	Chrysene-d12	14.01

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecanoic acid, butyl ester	340	C22H44O2	000123-95-5	91
2		n-Butyl myristate	284	C18H36O2	000110-36-1	90
3		Cyclopentanecarboxylic acid, 2-a...	129	C6H11NO2	040482-05-1	46
4		Decane, 2,2,3-trimethyl-	184	C13H28	062338-09-4	43
5		1-[N-Aziridyl]heptene-3	139	C9H17N	1000255-07-6	38



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF011116\
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Sample : H1043-18
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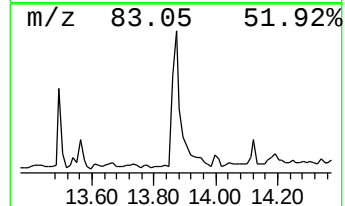
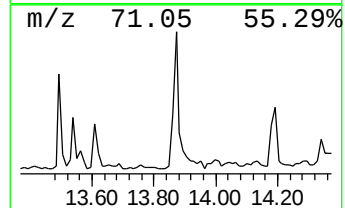
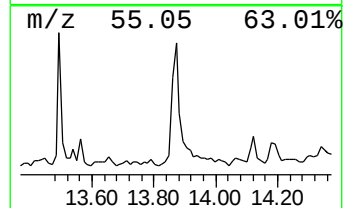
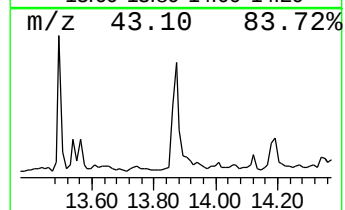
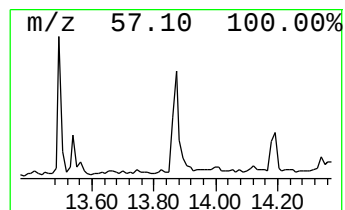
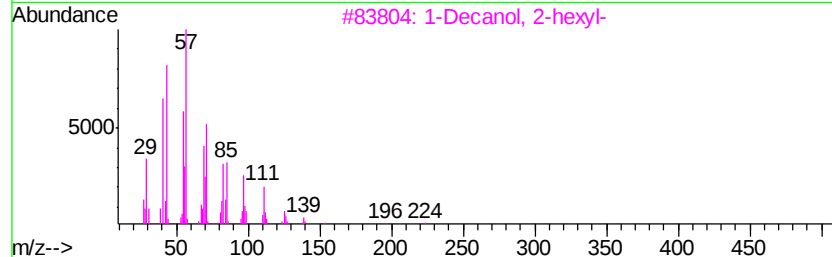
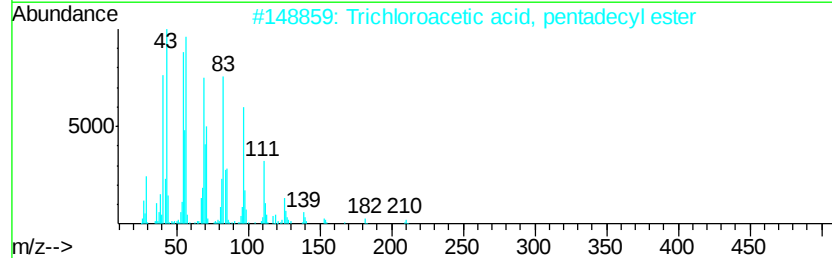
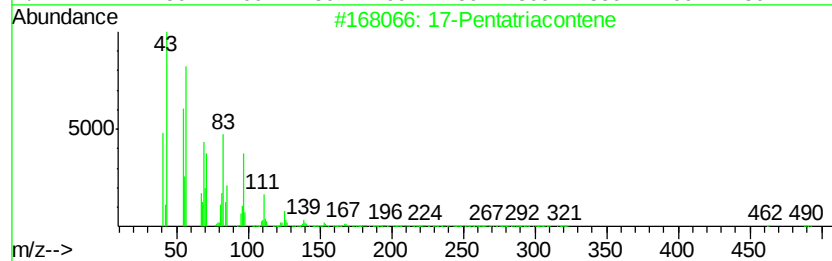
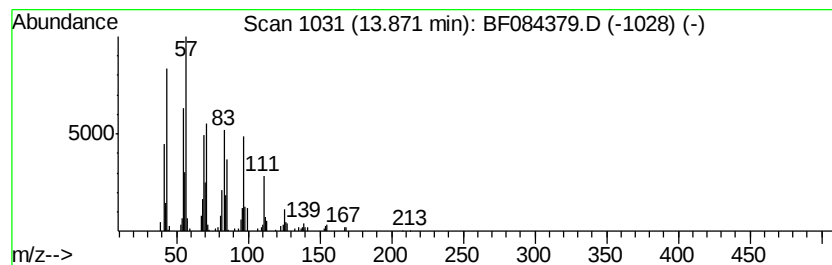
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 17-Pentatriacontene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.87	4.20 ng	267776	Chrysene-d12	14.01
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	17-Pentatriacontene		491 C35H70	006971-40-0 87
2	Trichloroacetic acid, pentadecyl...		372 C17H31Cl3O2	074339-53-0 76
3	1-Decanol, 2-hexyl-		242 C16H34O	002425-77-6 74
4	1-Docosene		308 C22H44	001599-67-3 68
5	2-Hexyl-1-octanol		214 C14H30O	019780-79-1 68



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF011116\
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Misc :
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Instrument :
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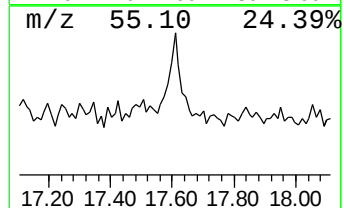
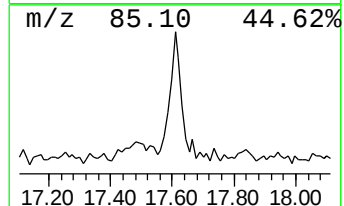
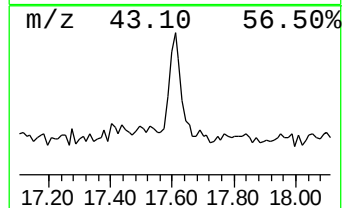
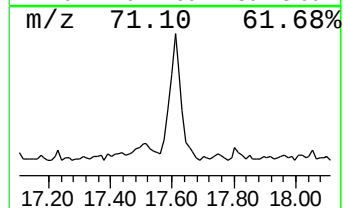
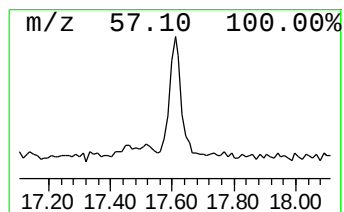
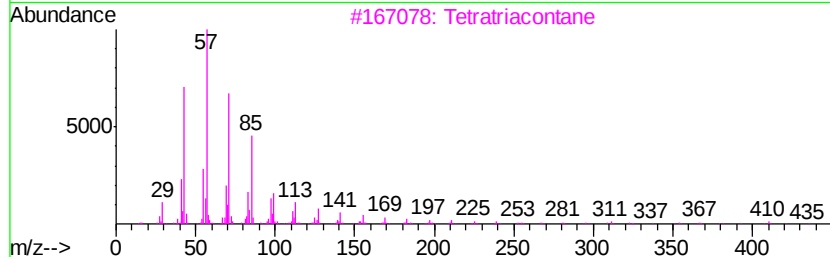
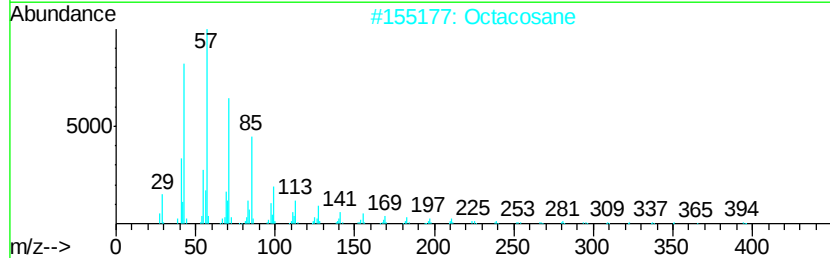
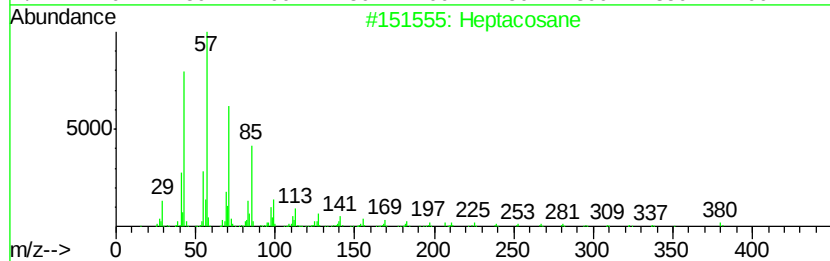
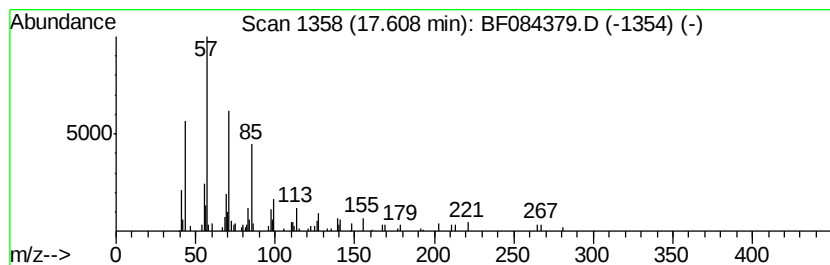
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Heptacosane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.61	2.27 ng	128958	Perylene-d12	15.45

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptacosane	380	C27H56	000593-49-7	91
2		Octacosane	394	C28H58	000630-02-4	90
3		Tetratriacontane	479	C34H70	014167-59-0	90
4		Hexatriacontane	507	C36H74	000630-06-8	87
5		Tetratetracontane	619	C44H90	007098-22-8	87



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF011116\
Data File : BF084379.D
Acq On : 11 Jan 2016 14:58
Operator : SJ/UM
Sample : H1043-18
Misc :
ALS Vial : 9 Sample Multiplier: 1

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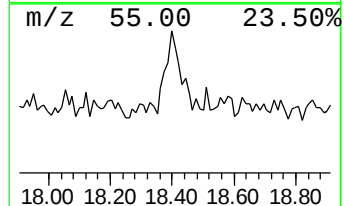
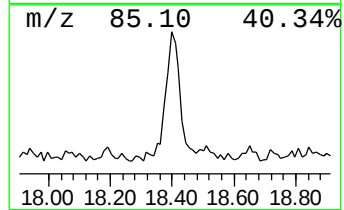
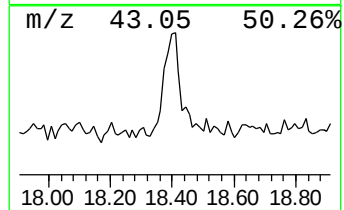
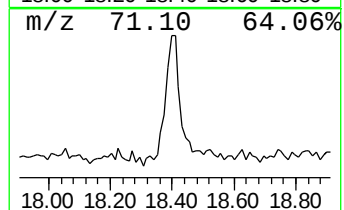
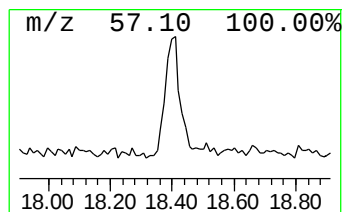
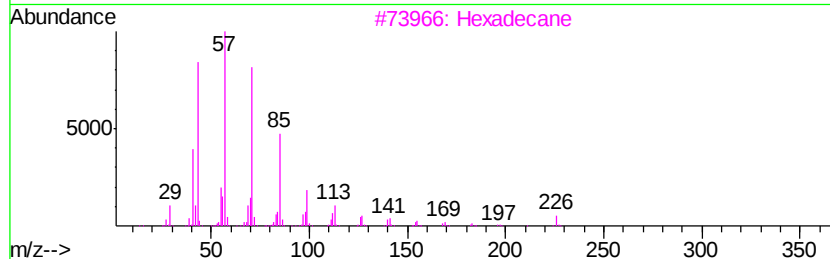
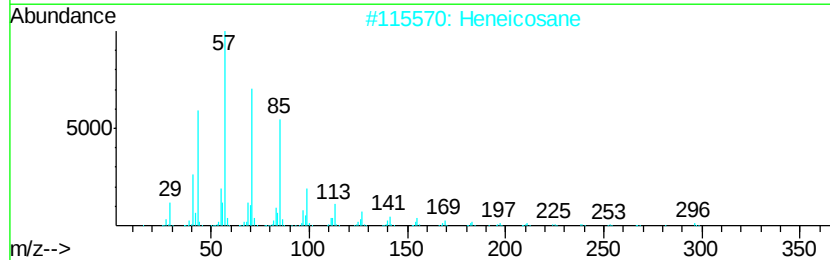
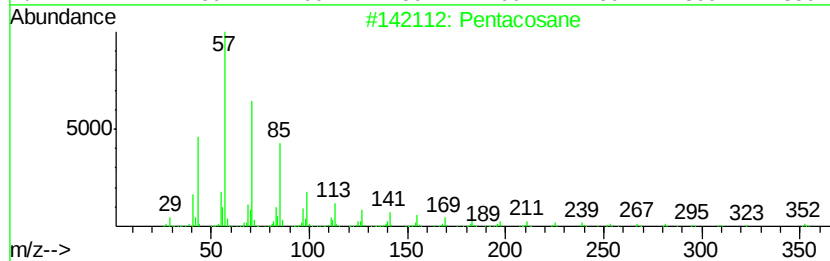
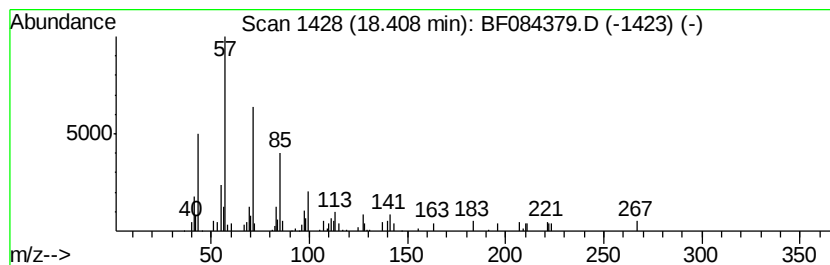
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Pentacosane Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.41	2.16 ng	122727	Perylene-d12	15.45

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentacosane	352	C25H52	000629-99-2	87
2		Heneicosane	296	C21H44	000629-94-7	83
3		Hexadecane	226	C16H34	000544-76-3	80
4		Hentriacontane	437	C31H64	000630-04-6	80
5		Triacontane	422	C30H62	000638-68-6	80



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF011116\
Data File : BF084379.D
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Misc :
ALS Vial : 9 Sample Multiplier: 1

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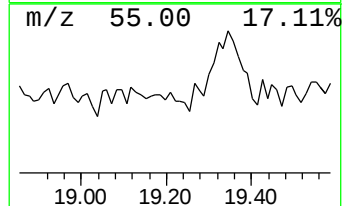
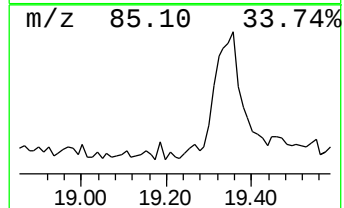
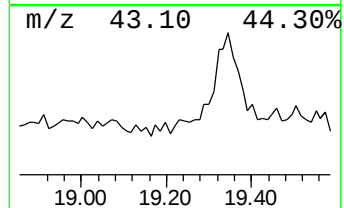
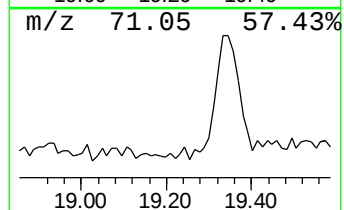
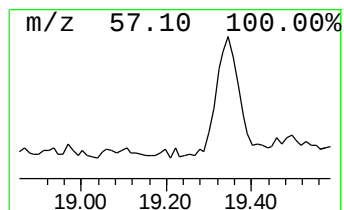
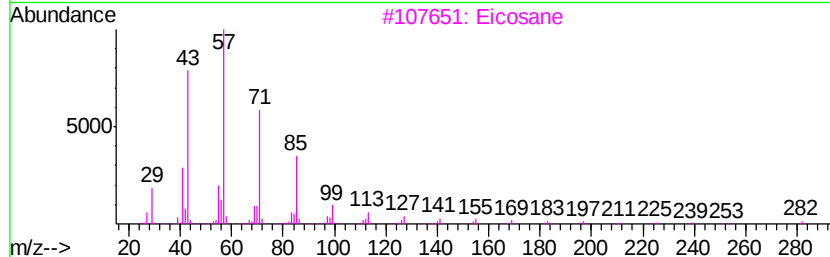
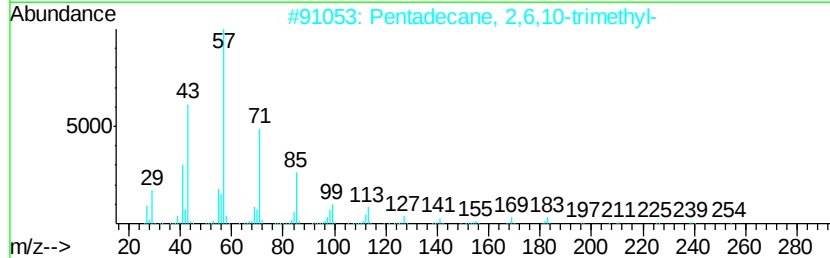
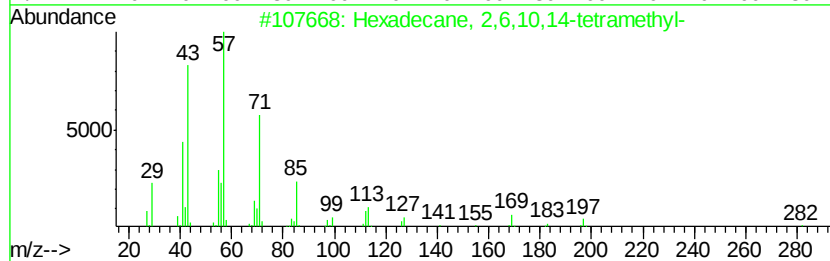
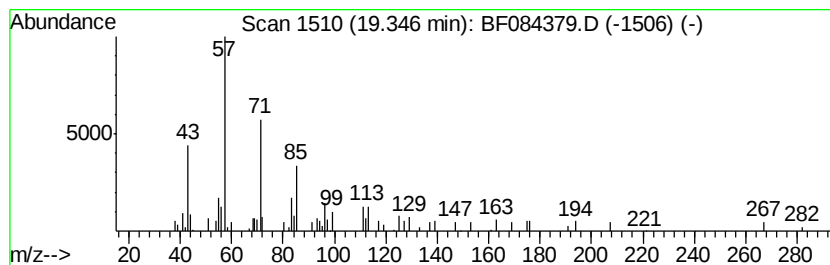
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Hexadecane, 2,6,10,14-tetra... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.35	2.25 ng	127618	Perylene-d12	15.45

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecane, 2,6,10,14-tetramethyl-	282	C20H42	000638-36-8	68
2		Pentadecane, 2,6,10-trimethyl-	254	C18H38	003892-00-0	64
3		Eicosane	282	C20H42	000112-95-8	62
4		Octadecane	254	C18H38	000593-45-3	59
5		1-Bromoeicosane	360	C20H41Br	004276-49-7	59



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF011116\
Data File : BF084379.D
Acq On : 11 Jan 2016 14:58
Operator : SJ/UM
Sample : H1043-18
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-4COMP

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF123115.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...	5.05	27.5	ng	2184010	1	6.85	1587020	20.0
unknown6.62	6.62	97.9	ng	7767760	1	6.85	1587020	20.0
Hexadecane	9.82	2.1	ng	213950	3	9.89	2080340	20.0
Pentadecane	10.54	2.2	ng	231934	3	9.89	2080340	20.0
Benzenesulfonamid...	10.80	8.7	ng	815253	4	11.38	1883230	20.0
Hexadecanoic acid...	12.79	4.9	ng	314888	5	14.01	1273880	20.0
Phenol, 4,4'-(1-m...	12.85	4.8	ng	304784	5	14.01	1273880	20.0
Octadecanoic acid...	13.49	3.5	ng	221604	5	14.01	1273880	20.0
17-Pentatriacontene	13.87	4.2	ng	267776	5	14.01	1273880	20.0
Heptacosane	17.61	2.3	ng	128958	6	15.45	1136760	20.0
Pentacosane	18.41	2.2	ng	122727	6	15.45	1136760	20.0
Hexadecane, 2,6,1...	19.35	2.3	ng	127618	6	15.45	1136760	20.0