

Data Path : Z:\HPCHEM1\BNA F\DATA\BF011016\
 Data File : BF084366.D
 Acq On : 10 Jan 2016 15:52
 Operator : SJ/UM
 Sample : H1043-16
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-3COMP

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	257	259	269	rVB	1437819	2351655	26.60%	3.059%
2	5.482	294	297	299	rBV	7060715	8795024	99.47%	11.440%
3	6.510	384	387	389	rBV	7138965	8034722	90.87%	10.451%
4	6.636	395	398	400	rBV	7663842	8841864	100.00%	11.501%
5	6.853	415	417	419	rBV	2063111	1660622	18.78%	2.160%
6	6.922	421	423	427	rBV	133243	118433	1.34%	0.154%
7	7.013	428	431	433	rVB	6831520	6096089	68.95%	7.929%
8	7.425	463	467	469	rBV	5193460	5450148	61.64%	7.089%
9	8.145	527	530	532	rBV	2179486	2190019	24.77%	2.849%
10	9.219	621	624	627	rBV	7487040	7525308	85.11%	9.788%
11	9.493	646	648	650	rBV	421835	334813	3.79%	0.435%
12	9.619	656	659	660	rBV	1268801	1433918	16.22%	1.865%
13	9.825	676	677	681	rBV2	148062	237013	2.68%	0.308%
14	9.893	681	683	686	rVB	2536190	2303639	26.05%	2.996%
15	10.328	719	721	723	rVB	435778	388852	4.40%	0.506%
16	10.545	737	740	744	rBV	123837	246298	2.79%	0.320%
17	10.682	749	752	755	rVV2	4024370	5750743	65.04%	7.480%
18	10.796	759	762	770	rVB2	624469	974399	11.02%	1.267%
19	10.991	777	779	784	rBV3	48419	120980	1.37%	0.157%
20	11.254	800	802	803	rBV	241253	218243	2.47%	0.284%
21	11.379	811	813	815	rBV	2648660	2257006	25.53%	2.936%
22	11.436	815	818	822	rVB	45292	96383	1.09%	0.125%
23	12.797	935	937	939	rBV	362505	391681	4.43%	0.509%
24	12.854	939	942	946	rVV2	187094	347355	3.93%	0.452%
25	12.968	949	952	955	rBV	6340648	6673101	75.47%	8.680%
26	13.494	996	998	1000	rBV	283514	288444	3.26%	0.375%
27	13.871	1028	1031	1041	rBV	260079	399798	4.52%	0.520%
28	14.008	1041	1043	1046	rBV	1114331	1515265	17.14%	1.971%
29	14.340	1070	1072	1074	rBV	122531	146259	1.65%	0.190%
30	14.797	1108	1112	1114	rVB2	56494	95803	1.08%	0.125%
31	14.865	1115	1118	1120	rBV	229137	220682	2.50%	0.287%
32	15.437	1165	1168	1171	rBV	824059	1280500	14.48%	1.666%
33	16.385	1248	1251	1257	rVB2	41872	95382	1.08%	0.124%

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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-BF123115.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

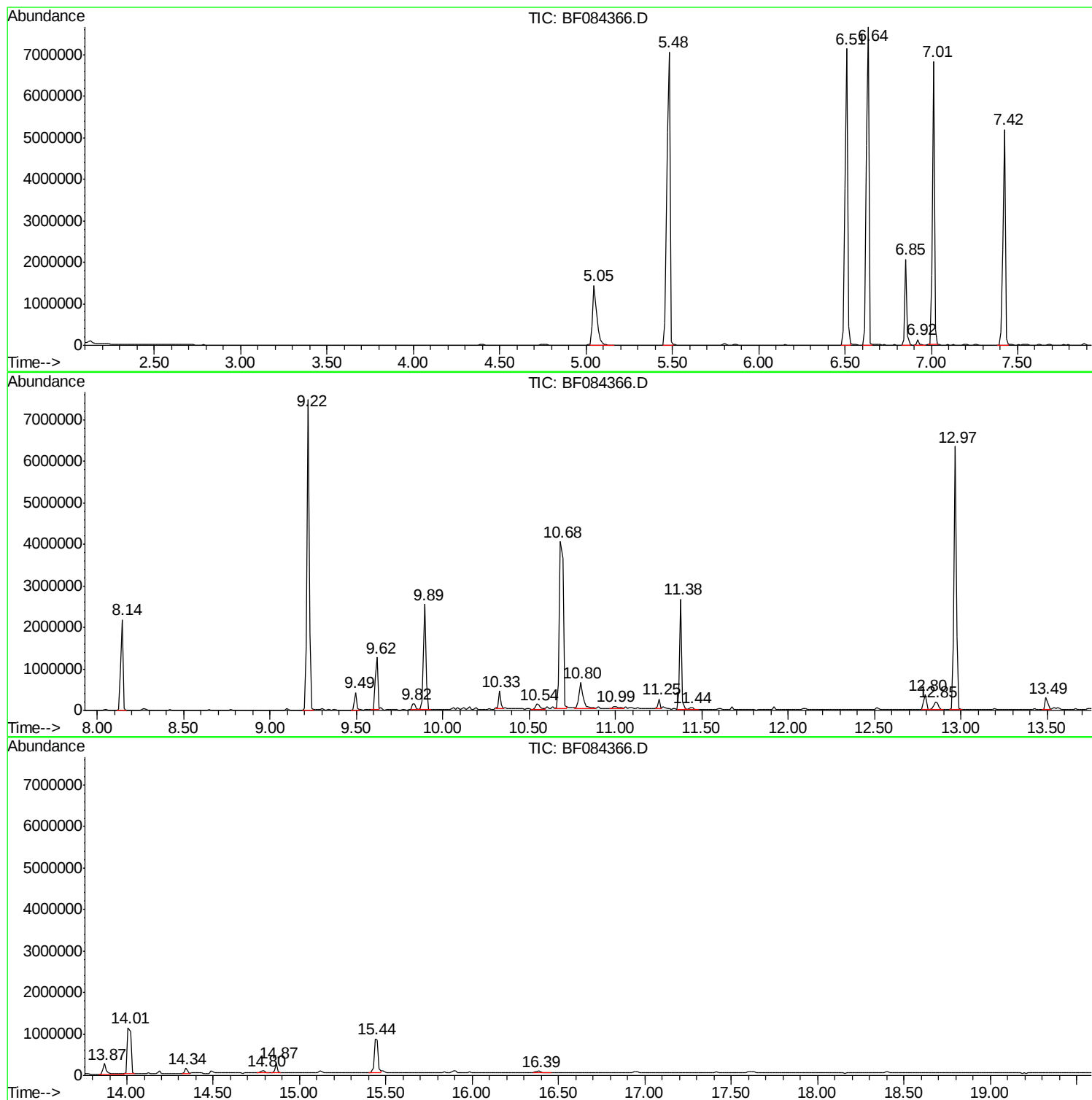
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF123115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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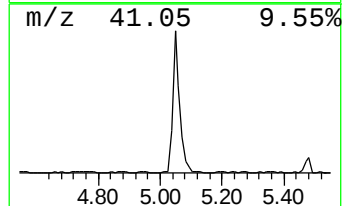
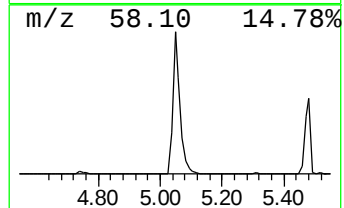
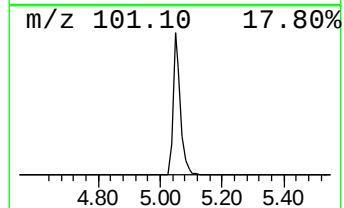
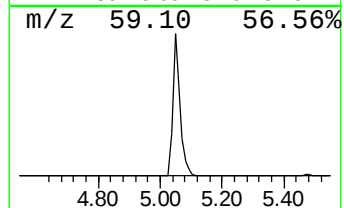
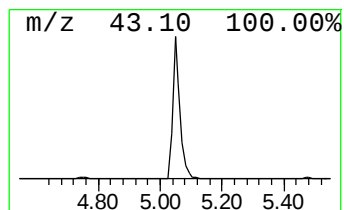
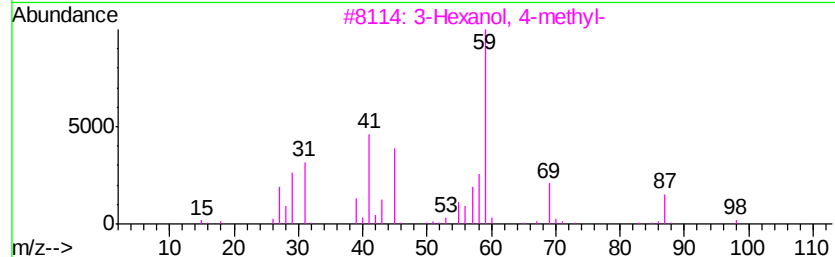
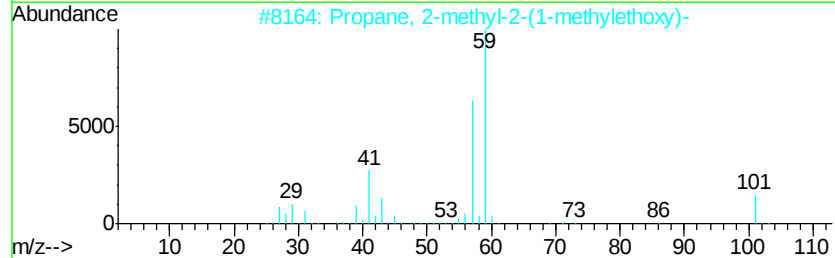
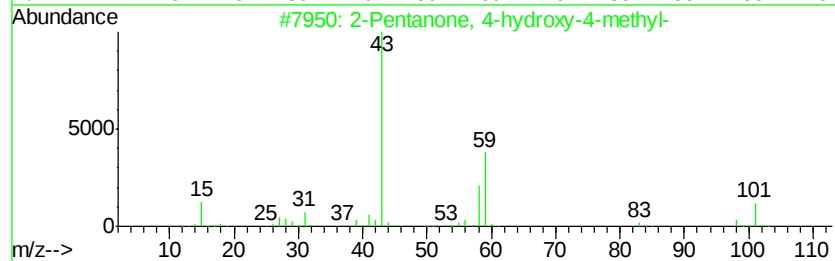
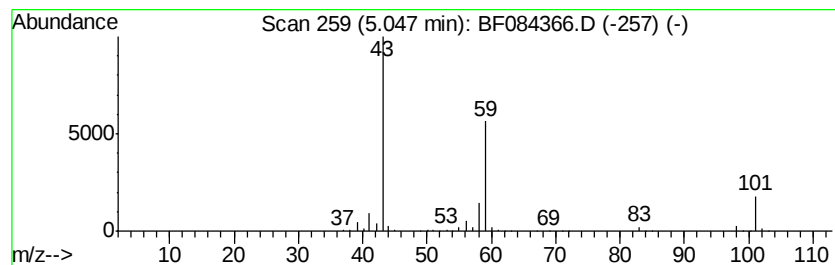
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.05	28.32 ng	2351660	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	59
2			Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	36
3			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
4			Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	32
5			1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10



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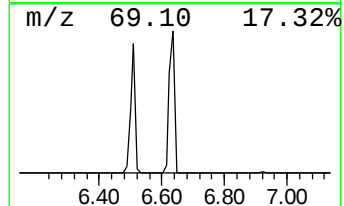
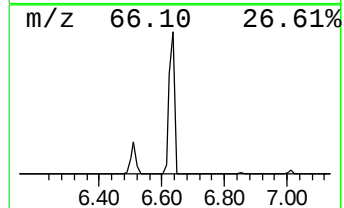
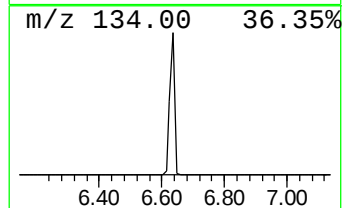
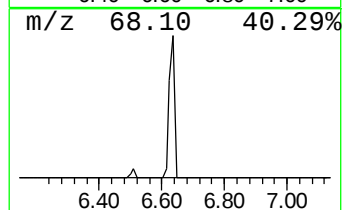
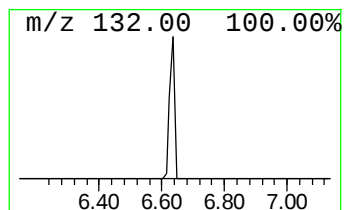
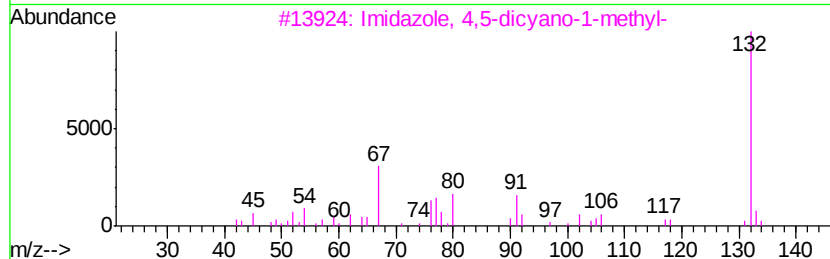
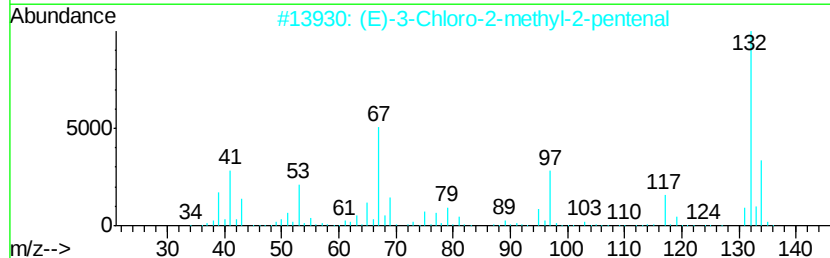
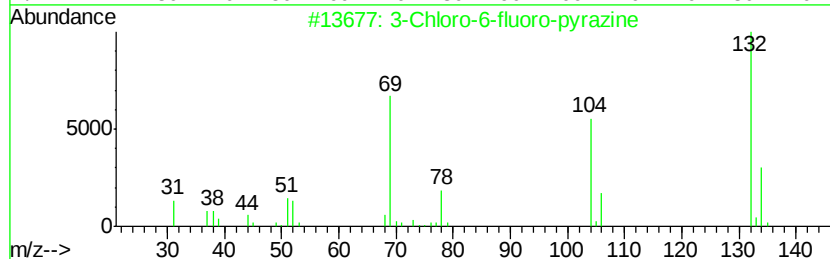
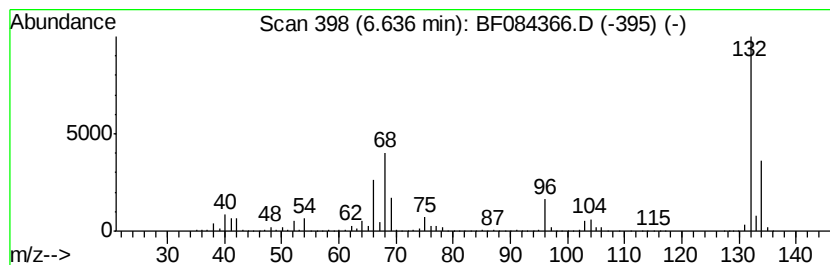
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 2 unknown6.64 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.64	106.49 ng	8841860	1,4-Dichlorobenzene-d4	6.85

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	16
3		Imidazole, 4,5-dicyano-1-methyl-	132	C6H4N4	019485-35-9	9
4		Benzene, 1,2,4-trifluoro-	132	C6H3F3	000367-23-7	9
5		Benzene, 1,3,5-trifluoro-	132	C6H3F3	000372-38-3	9



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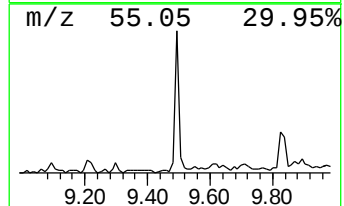
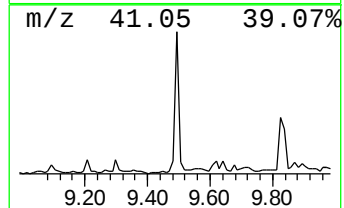
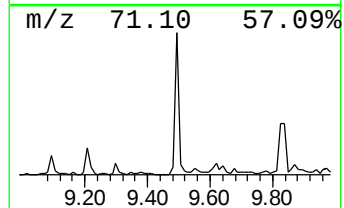
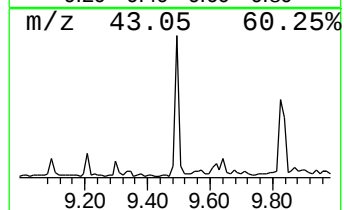
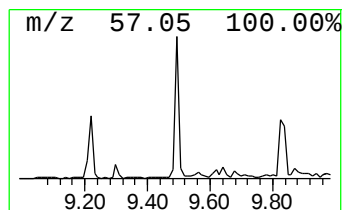
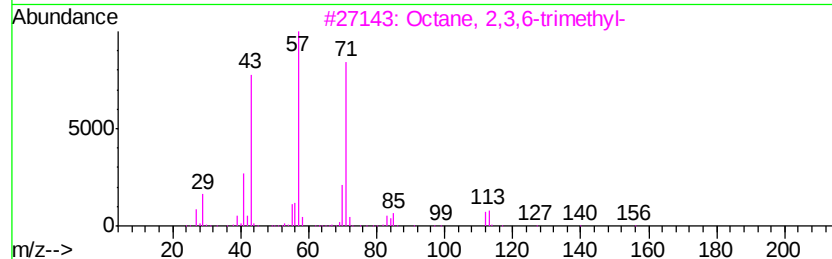
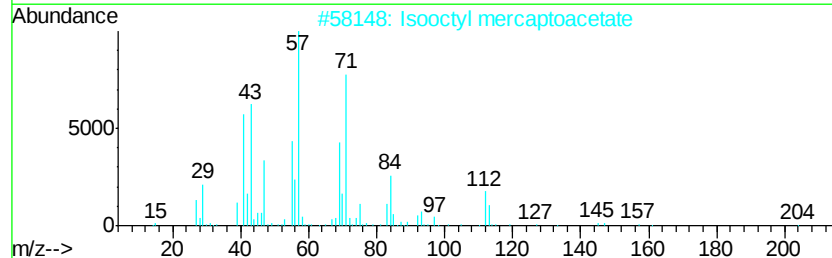
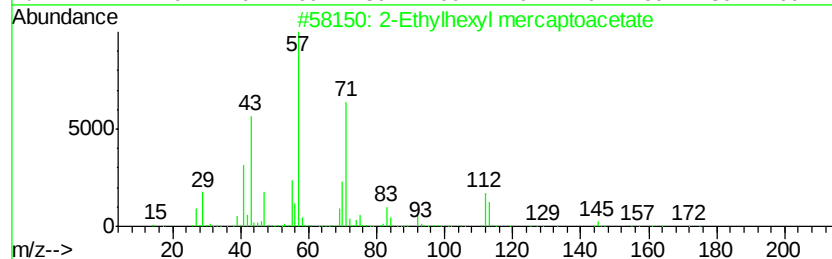
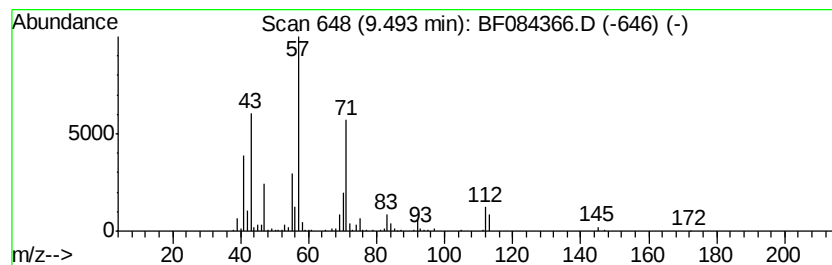
Instrument :
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 2-Ethylhexyl mercaptoacetate Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD		R.T.
9.49	2.91 ng	334813	Acenaphthene-d10		9.89
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Ethylhexyl mercaptoacetate	204	C10H20O2S	007659-86-1	80
2	Isooctyl mercaptoacetate	204	C10H20O2S	025103-09-7	59
3	Octane, 2,3,6-trimethyl-	156	C11H24	062016-33-5	59
4	Decane, 4-methyl-	156	C11H24	002847-72-5	53
5	Heptane, 3-[(ethenyloxy)methyl]-	156	C10H20O	000103-44-6	50



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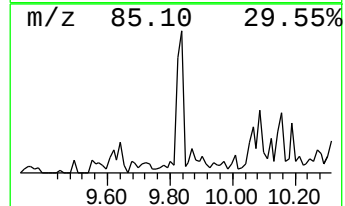
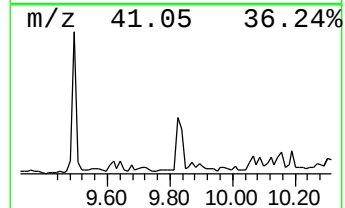
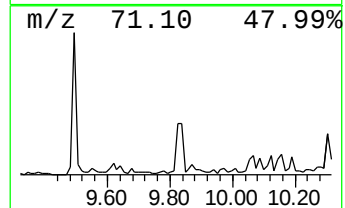
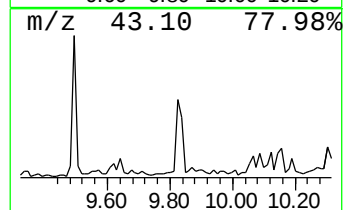
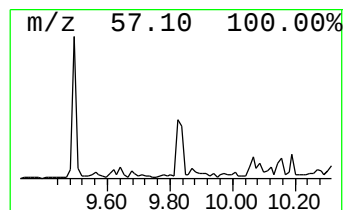
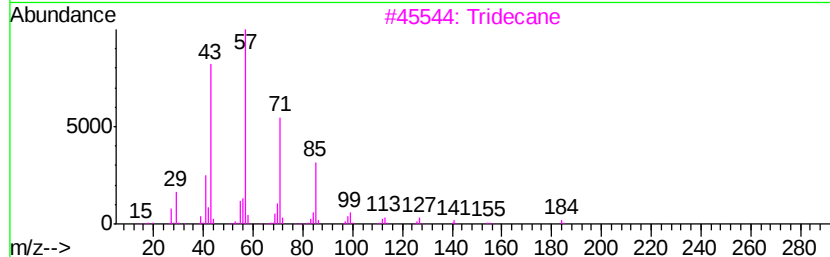
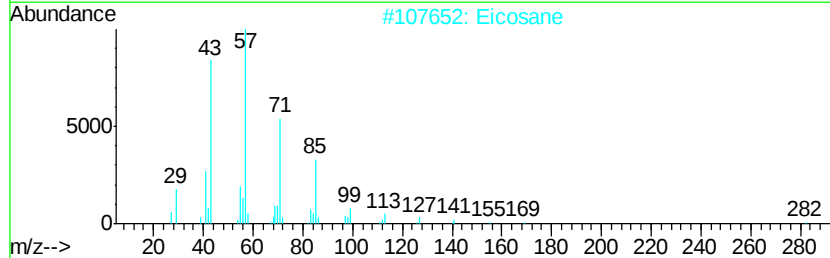
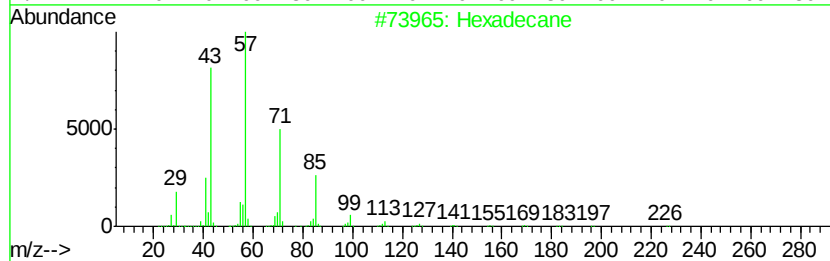
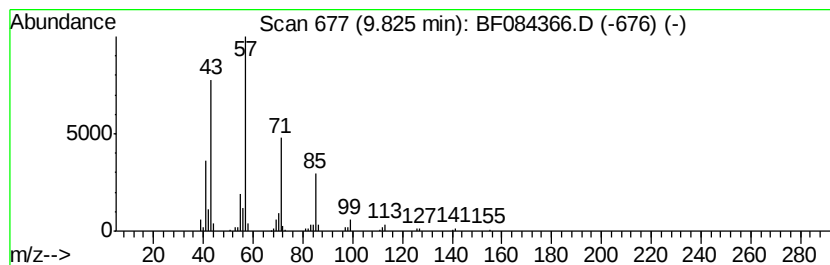
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Hexadecane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.82	2.06 ng	237013	Acenaphthene-d10	9.89

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane		226	C16H34	000544-76-3	91
2	Eicosane		282	C20H42	000112-95-8	86
3	Tridecane		184	C13H28	000629-50-5	78
4	Dodecane, 2-methyl-		184	C13H28	001560-97-0	78
5	Undecane, 2-methyl-		170	C12H26	007045-71-8	72



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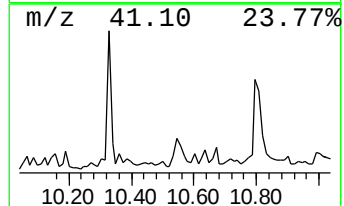
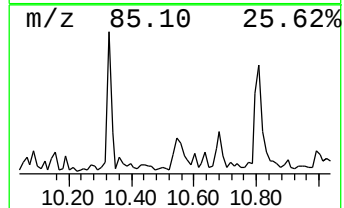
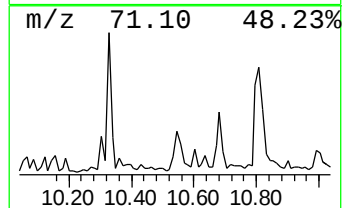
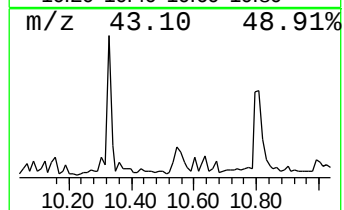
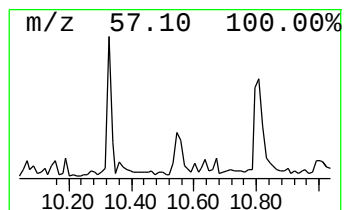
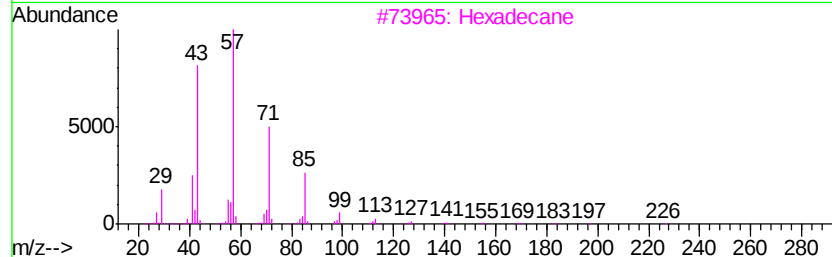
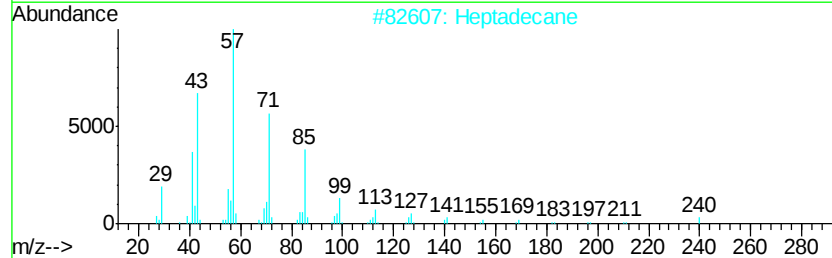
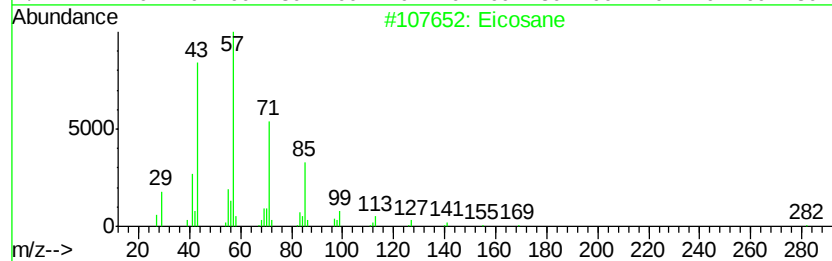
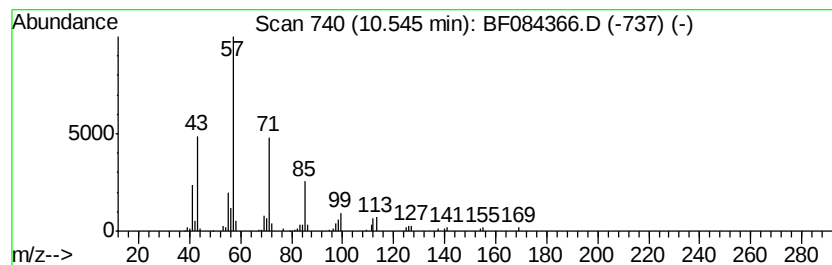
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 7 Eicosane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.54	2.14 ng	246298	Acenaphthene-d10	9.89

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	90
2		Heptadecane	240	C17H36	000629-78-7	80
3		Hexadecane	226	C16H34	000544-76-3	72
4		Hentriacontane	437	C31H64	000630-04-6	72
5		Pentadecane	212	C15H32	000629-62-9	72



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ALS Vial : 17 Sample Multiplier: 1

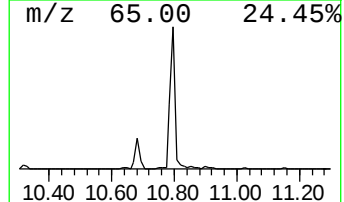
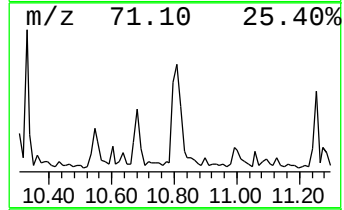
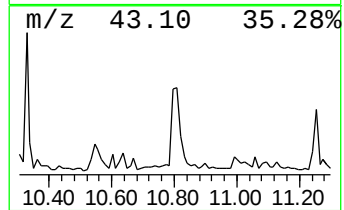
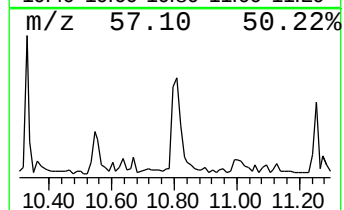
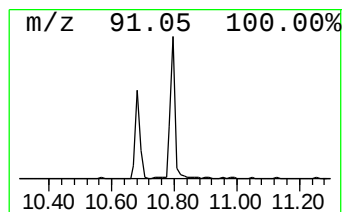
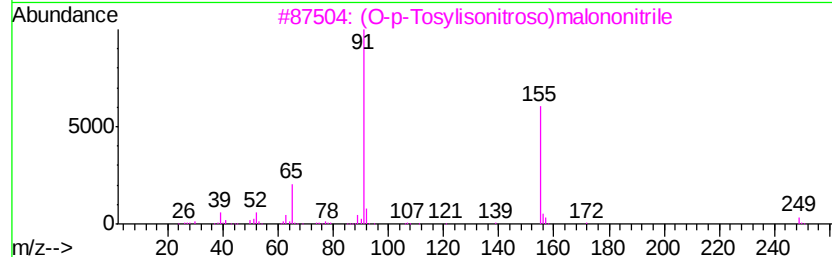
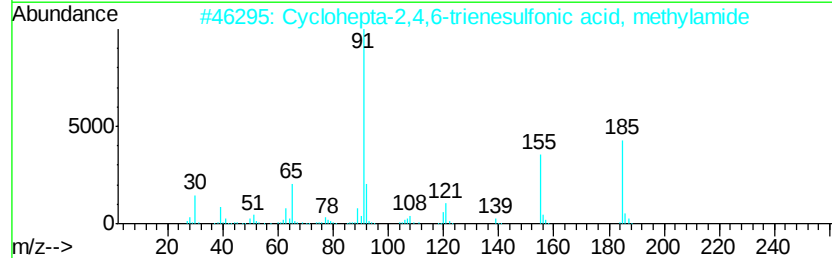
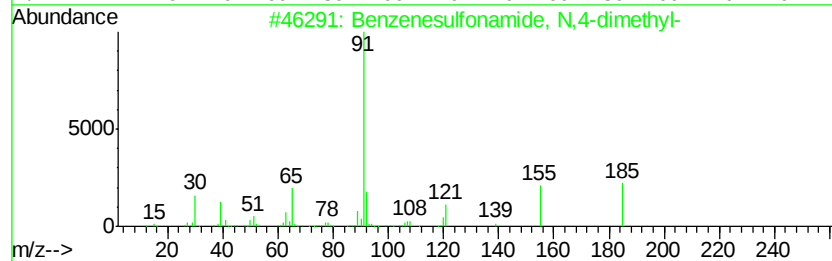
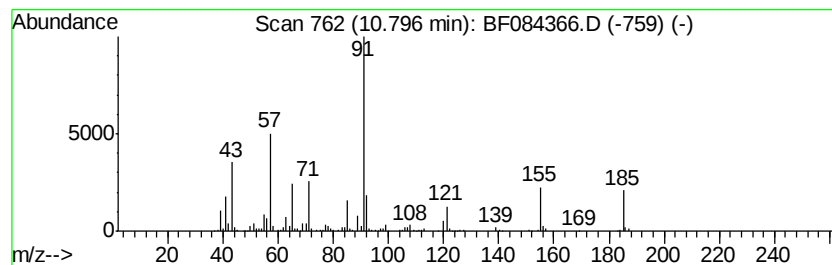
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Benzenesulfonamide, N,4-dim... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.80	8.63 ng	974399	Phenanthrene-d10	11.38	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzenesulfonamide, N,4-dimethyl-	185	C8H11NO2S	000640-61-9	89
2	Cyclohepta-2,4,6-trienesulfonic ...	185	C8H11NO2S	1000187-28-7	38
3	(O-p-Tosvlisonitroso)malononitrile	249	C10H7N3O3S	020893-01-0	38
4	4-Toluenesulfonylmethylisocyanide	195	C9H9NO2S	036635-61-7	38
5	Ethyl(E/Z)-.alpha.-cyano-4-nitro...	416	C19H16N2O7S	1000287-26-8	38



Data Path : Z:\HPCHEM1\BNA F\DATA\BF011016\
Data File : BF084366.D
Acq On : 10 Jan 2016 15:52
Operator : SJ/UM
Sample : H1043-16
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-3COMP

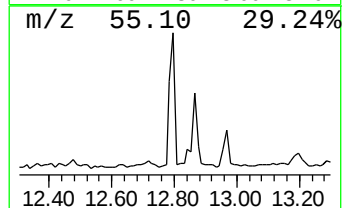
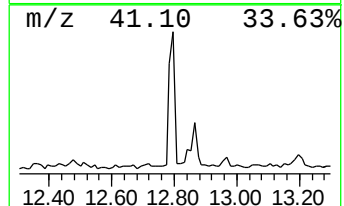
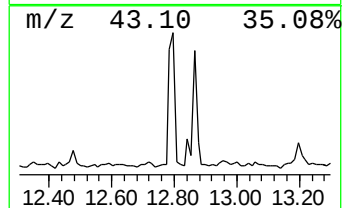
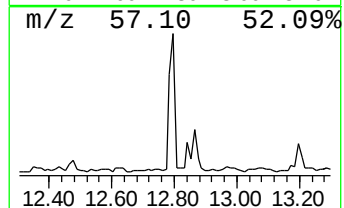
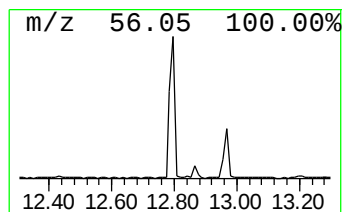
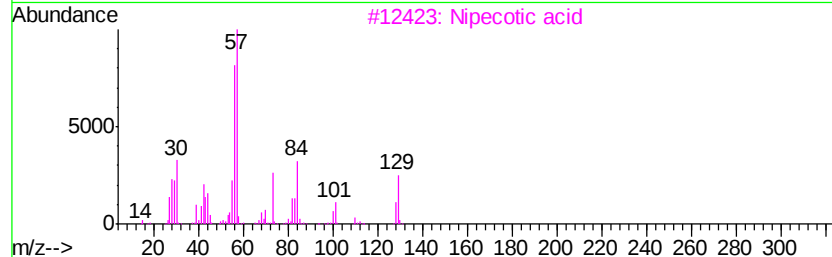
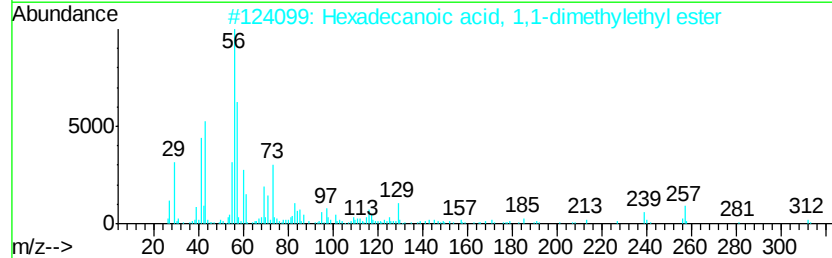
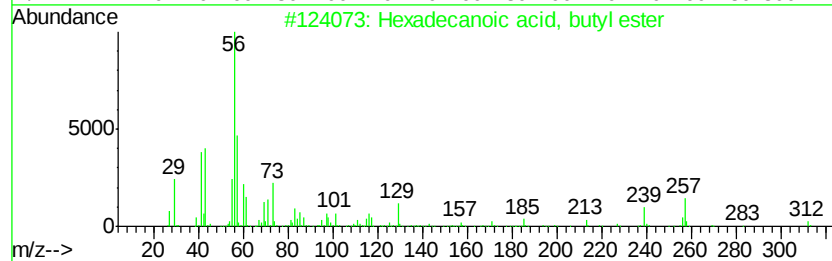
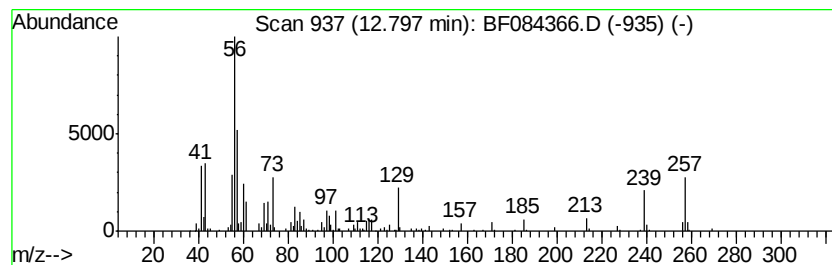
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Hexadecanoic acid, butyl ester Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.80	5.17 ng	391681	Chrysene-d12	14.02

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecanoic acid, butyl ester	312	C20H40O2	000111-06-8	81
2		Hexadecanoic acid, 1,1-dimethyle...	312	C20H40O2	031158-91-5	68
3		Nipecotic acid	129	C6H11NO2	000498-95-3	47
4		Cyclobutane, 1,1-dimethyl-2-octyl-	196	C14H28	062338-30-1	35
5		Hexadecanoic acid, 2-methylpropy...	312	C20H40O2	000110-34-9	32



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF011016\
Data File : BF084366.D
Acq On : 10 Jan 2016 15:52
Operator : SJ/UM
Sample : H1043-16
Misc :
ALS Vial : 17 Sample Multiplier: 1

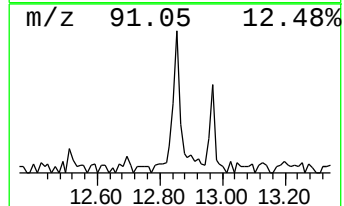
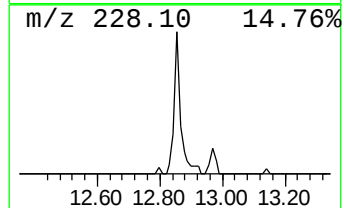
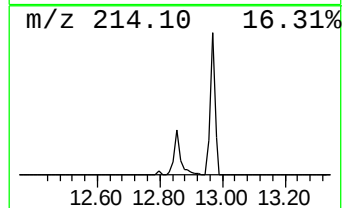
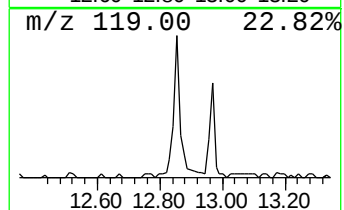
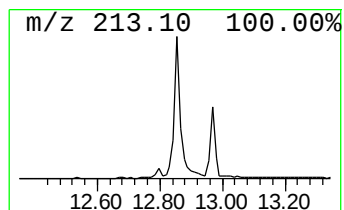
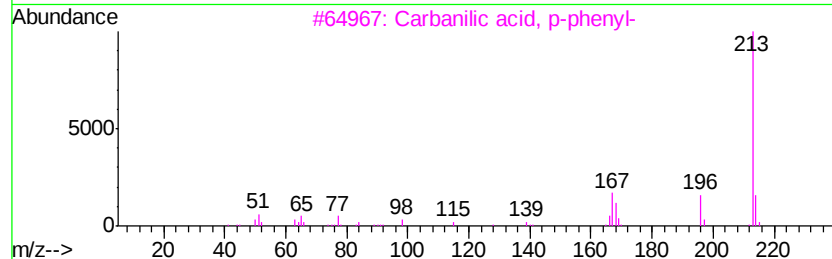
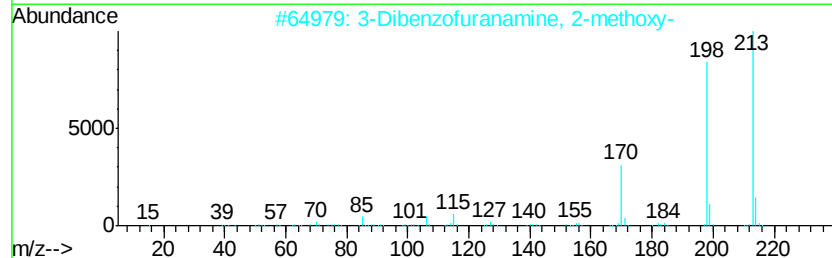
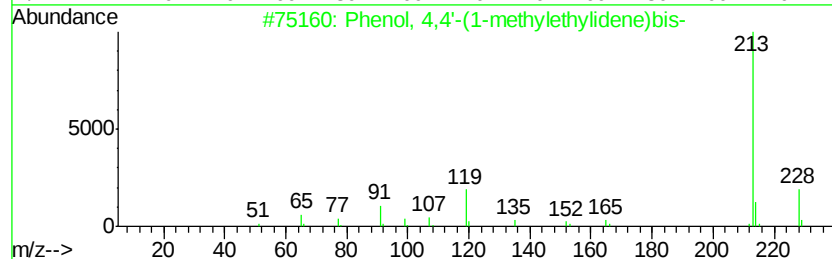
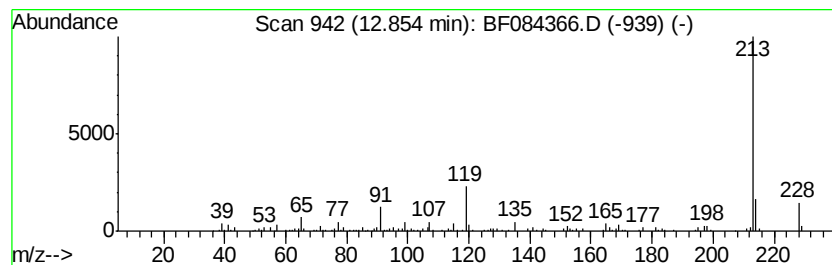
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF123115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Phenol, 4,4'-(1-methylethyl... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.85	4.58 ng	347355	Chrysene-d12	14.02	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 4,4'-(1-methylethylidene...	228	C15H16O2	000080-05-7	96
2	3-Dibenzofuranamine, 2-methoxy-	213	C13H11NO2	005834-17-3	47
3	Carbanilic acid, p-phenyl-	213	C13H11NO2	004474-53-7	47
4	2,4(1H,3H)-Pyrimidinedione, 1,3-...	228	C9H16N2O3Si	089447-84-7	42
5	Phenol, 2,4'-isopropylidenedi-	228	C15H16O2	000837-08-1	41



Data Path : Z:\HPCHEM1\BNA F\DATA\BF011016\
Data File : BF084366.D
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Misc :
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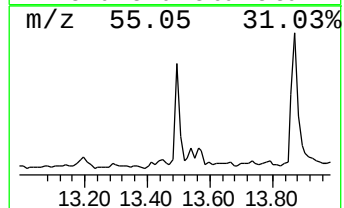
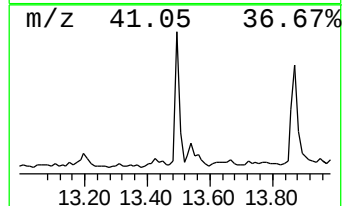
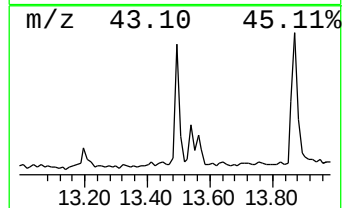
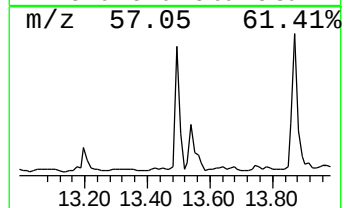
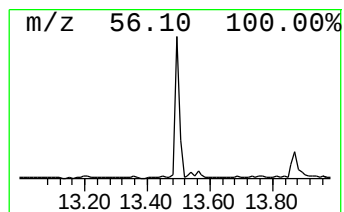
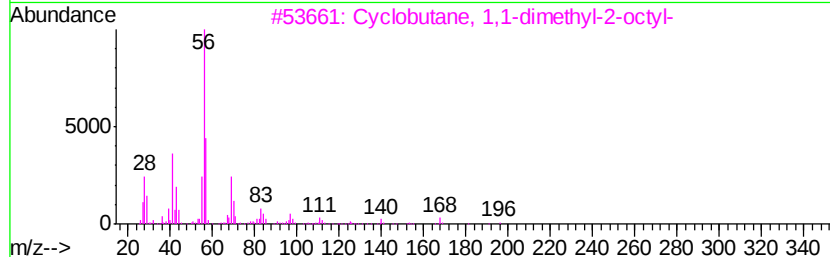
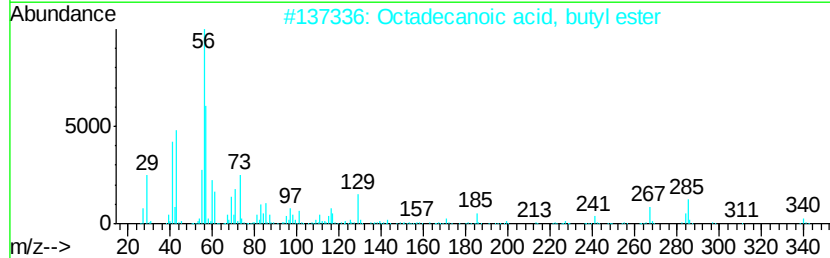
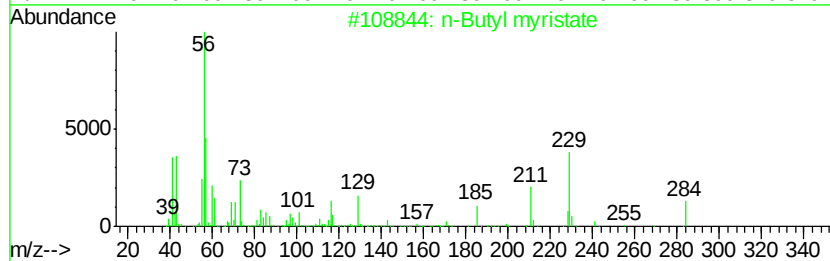
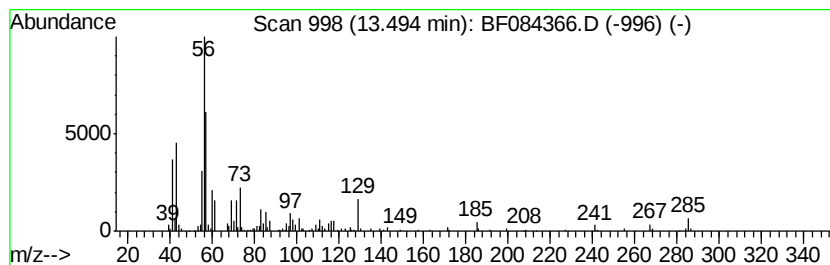
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 n-Butyl myristate Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.49	3.81 ng	288444	Chrysene-d12	14.02

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Butyl myristate	284	C18H36O2	000110-36-1	90
2			Octadecanoic acid, butyl ester	340	C22H44O2	000123-95-5	83
3			Cyclobutane, 1,1-dimethyl-2-octyl-	196	C14H28	062338-30-1	43
4			1-Undecene, 2-methyl-	168	C12H24	018516-37-5	38
5			Cyclobutane, ethyl-	84	C6H12	004806-61-5	38



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF011016\
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Sample : H1043-16
Misc :
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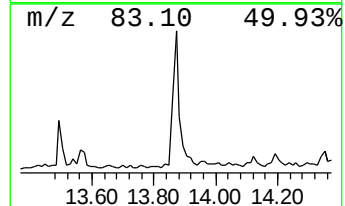
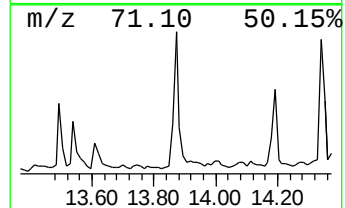
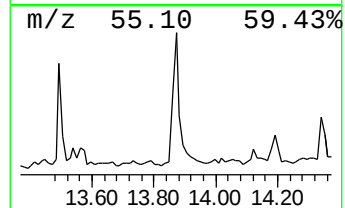
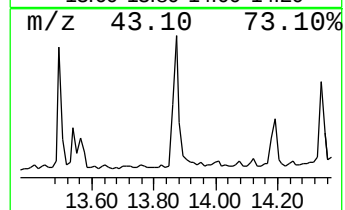
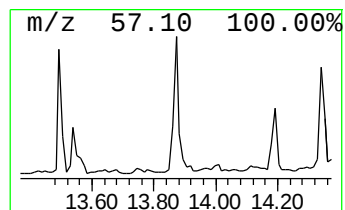
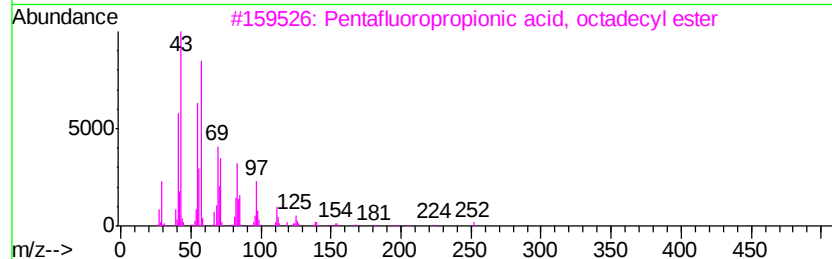
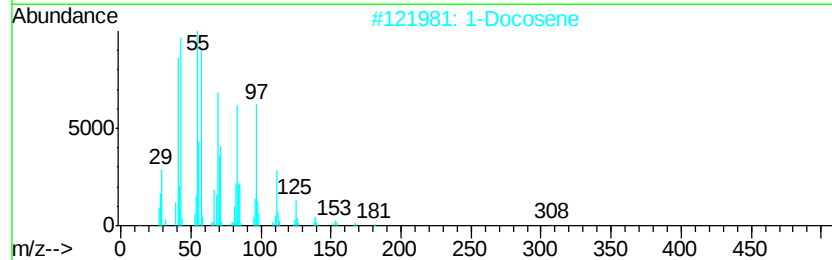
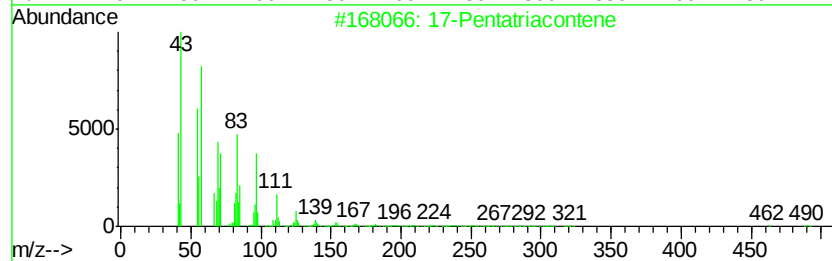
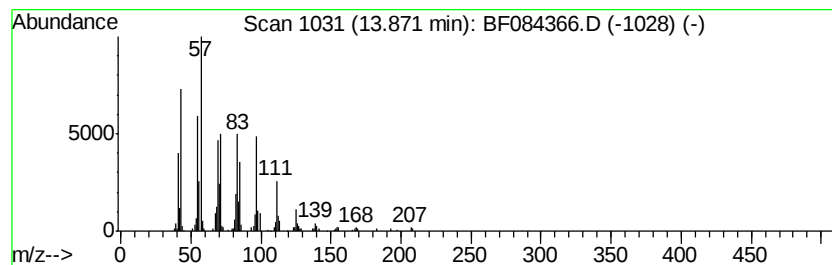
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 17-Pentatriacontene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.87	5.28 ng	399798	Chrysene-d12	14.02	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	17-Pentatriacontene	491	C35H70	006971-40-0	87
2	1-Docosene	308	C22H44	001599-67-3	76
3	Pentafluoropropionic acid, octadecyl...	416	C21H37F5O2	1000280-07-7	72
4	3-Chloropropionic acid, heptadecyl...	346	C20H39ClO2	1000283-05-1	62
5	Trichloroacetic acid, pentadecyl...	372	C17H31Cl3O2	074339-53-0	58



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF011016\
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ClientSampleId :
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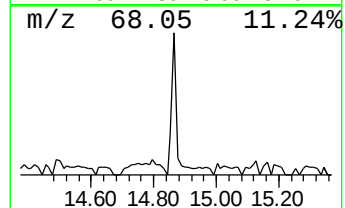
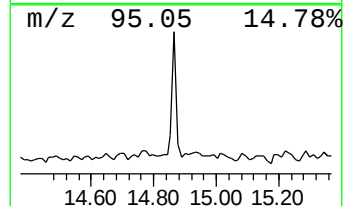
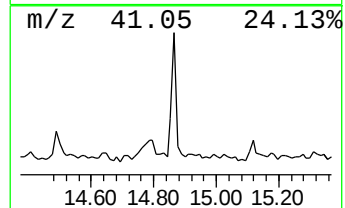
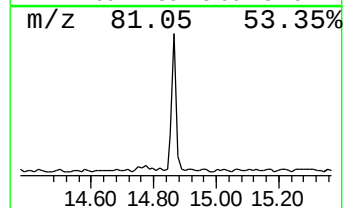
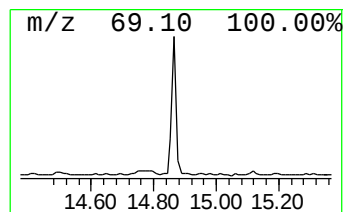
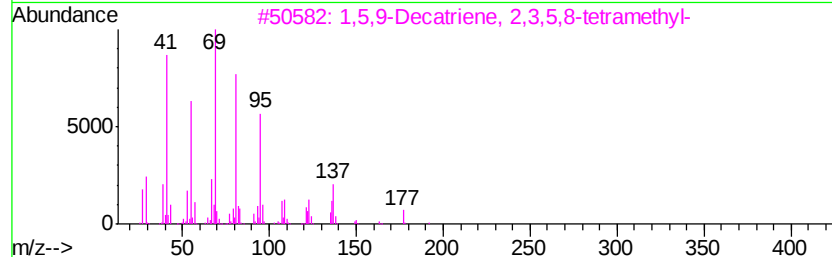
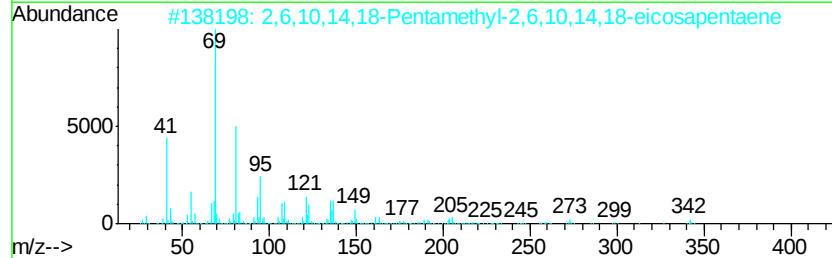
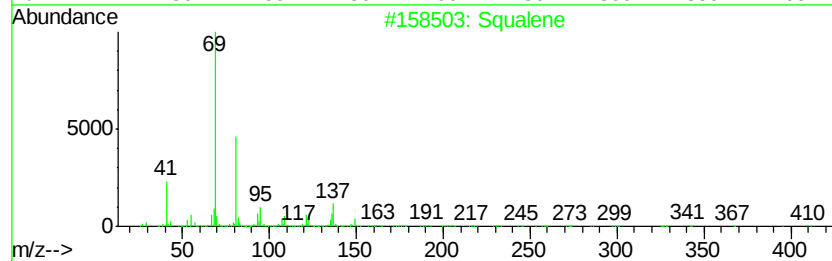
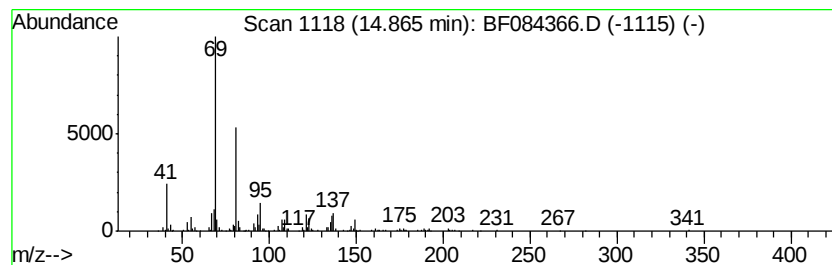
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Squalene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.87	3.45 ng	220682	Perylene-d12	15.45

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Squalene	410	C30H50	007683-64-9	91
2			2,6,10,14,18-Pentamethyl-2,6,10,...	342	C25H42	075581-03-2	83
3			1,5,9-Decatriene, 2,3,5,8-tetram...	192	C14H24	230646-72-7	80
4			1,5,9-Undecatriene, 2,6,10-trime...	192	C14H24	062951-96-6	80
5			7,11-Dimethyldodeca-2,6,10-trien...	208	C14H24O	125529-10-4	72



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF011016\
Data File : BF084366.D
Acq On : 10 Jan 2016 15:52
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Sample : H1043-16
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
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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	28.3	ng	2351660	1	6.85	1660620	20.0
unknown6.64	6.64	106.5	ng	8841860	1	6.85	1660620	20.0
2-Ethylhexyl merc...	9.49	2.9	ng	334813	3	9.89	2303640	20.0
Hexadecane	9.82	2.1	ng	237013	3	9.89	2303640	20.0
Eicosane	10.54	2.1	ng	246298	3	9.89	2303640	20.0
Benzenesulfonamid...	10.80	8.6	ng	974399	4	11.38	2257010	20.0
Hexadecanoic acid...	12.80	5.2	ng	391681	5	14.02	1515270	20.0
Phenol, 4,4'-(1-m...	12.85	4.6	ng	347355	5	14.02	1515270	20.0
n-Butyl myristate	13.49	3.8	ng	288444	5	14.02	1515270	20.0
17-Pentatriacontene	13.87	5.3	ng	399798	5	14.02	1515270	20.0
Squalene	14.87	3.5	ng	220682	6	15.45	1280500	20.0