

Data Path : Z:\HPCHEM1\BNA F\DATA\BF110816\
 Data File : BF091085.D
 Acq On : 8 Nov 2016 16:43
 Operator : UM/SJ
 Sample : H5571-01
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 FORMER-POTABLE-WELL

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-BF110316.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	1.755	4	6	51	rVB	3704213	10041238	100.00%	18.646%
2	4.796	269	272	277	rBV	355980	576998	5.75%	1.071%
3	5.253	309	312	322	rBV	1438507	2256252	22.47%	4.190%
4	6.304	403	404	412	rVV	880177	1422601	14.17%	2.642%
5	6.430	412	415	417	rVV	4211665	4088365	40.72%	7.592%
6	6.659	433	435	439	rBV	1140317	1114441	11.10%	2.069%
7	6.819	446	449	451	rBV	4150491	4322751	43.05%	8.027%
8	6.933	456	459	465	rVB3	63782	163624	1.63%	0.304%
9	7.230	482	485	490	rBV	2979319	3165535	31.53%	5.878%
10	7.870	537	541	545	rBV5	39704	134228	1.34%	0.249%
11	7.950	545	548	550	rBV	1324910	1501863	14.96%	2.789%
12	8.156	564	566	570	rVV4	79923	128002	1.27%	0.238%
13	8.293	575	578	579	rBV	77278	115661	1.15%	0.215%
14	8.408	585	588	593	rVB6	46490	138864	1.38%	0.258%
15	8.979	636	638	640	rVV3	67245	136789	1.36%	0.254%
16	9.025	640	642	645	rVV	4948653	5735862	57.12%	10.651%
17	9.070	645	646	647	rVV	138146	109846	1.09%	0.204%
18	9.425	676	677	679	rVV2	106992	117100	1.17%	0.217%
19	9.608	691	693	695	rVV2	125227	209030	2.08%	0.388%
20	9.699	697	701	703	rBV	1830496	1972867	19.65%	3.664%
21	9.825	710	712	718	rVB4	164522	241608	2.41%	0.449%
22	9.916	718	720	722	rBV2	68702	136486	1.36%	0.253%
23	10.019	727	729	731	rBV3	66640	135918	1.35%	0.252%
24	10.191	742	744	745	rBV	106170	120634	1.20%	0.224%
25	10.259	748	750	754	rVB3	83935	155521	1.55%	0.289%
26	10.362	757	759	762	rBV2	126198	249746	2.49%	0.464%
27	10.488	767	770	773	rVV	3116946	3271127	32.58%	6.074%
28	10.533	773	774	776	rVV2	128420	161540	1.61%	0.300%
29	10.625	780	782	785	rVB	230973	334816	3.33%	0.622%
30	11.093	821	823	825	rVB	194562	245201	2.44%	0.455%
31	11.174	828	830	833	rBV	1638995	1558526	15.52%	2.894%
32	11.642	869	871	874	rBV3	83868	151097	1.50%	0.281%
33	11.722	877	878	882	rVB3	140522	224966	2.24%	0.418%
34	12.762	967	969	972	rBV	4845306	6110187	60.85%	11.346%

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Sampling : 1 Min Area: 3 % of largest Peak
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Stop Thrs : 0 Peak Location: TOP

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35	13.665	1045	1048	1058	rVB	173874	246815	2.46%	0.458%
36	13.802	1058	1060	1067	rVB	1741712	1681908	16.75%	3.123%
37	14.648	1132	1134	1136	rBV	232574	210481	2.10%	0.391%
38	15.174	1178	1180	1183	rBV	804571	1162667	11.58%	2.159%

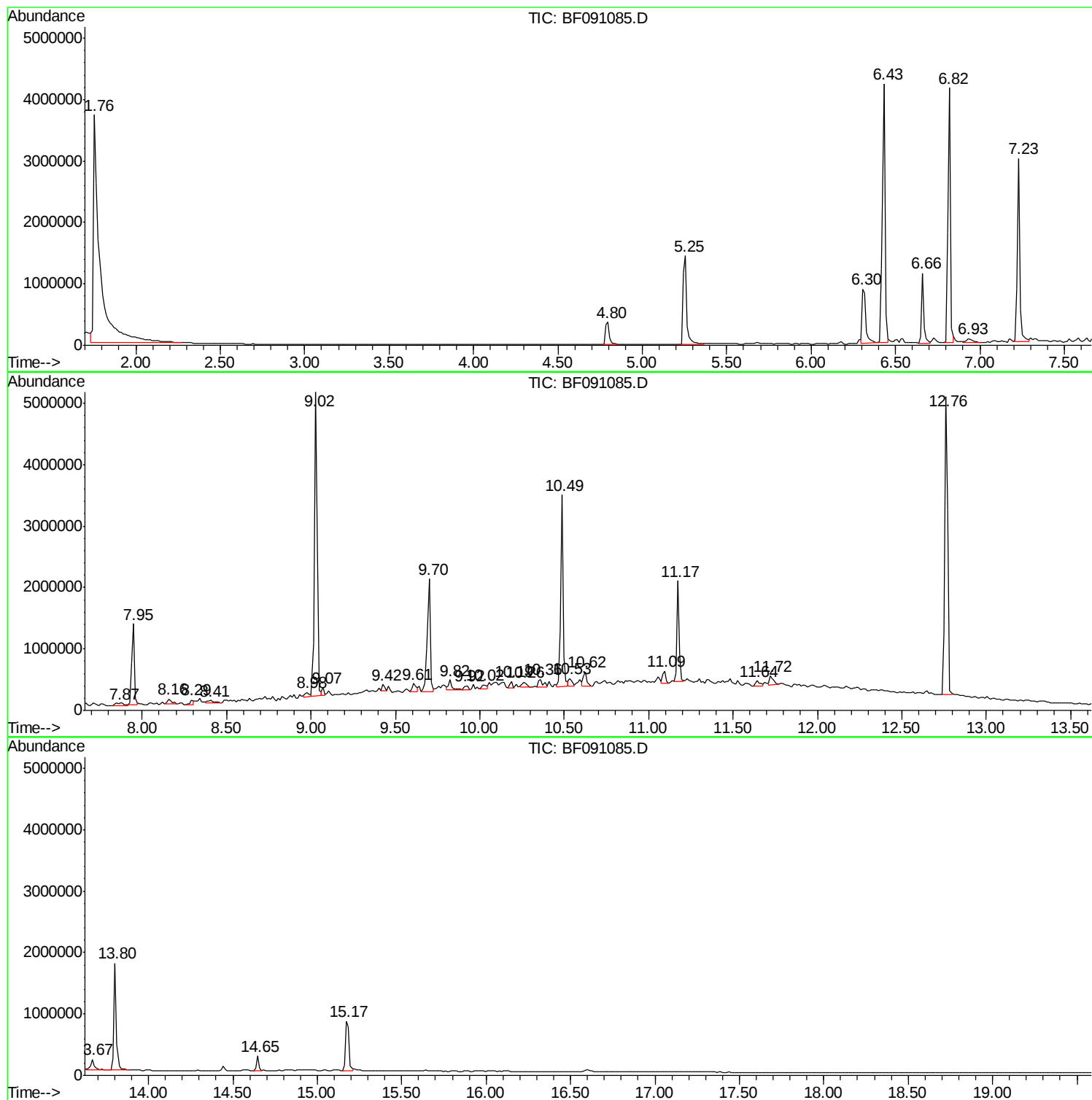
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF110316.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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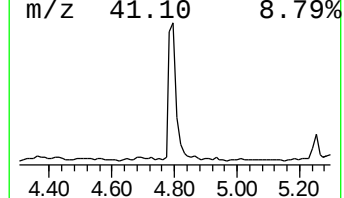
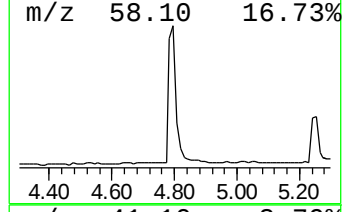
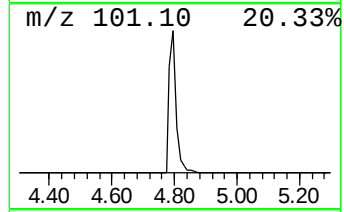
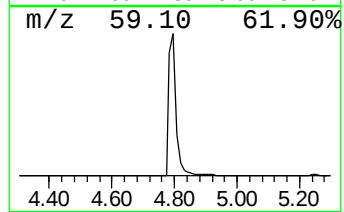
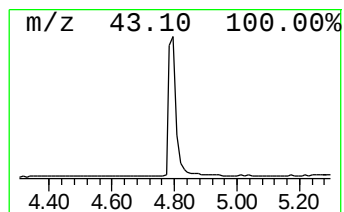
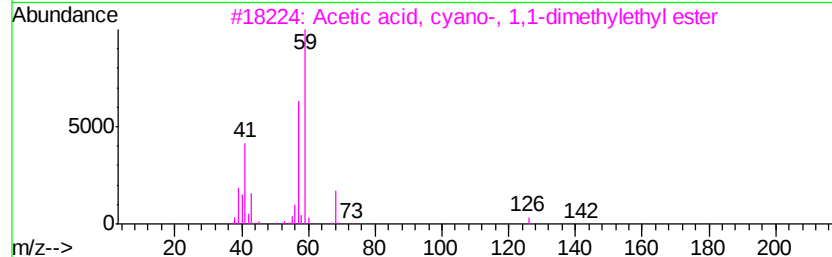
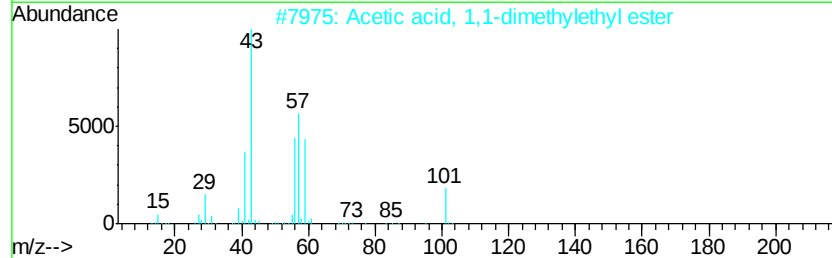
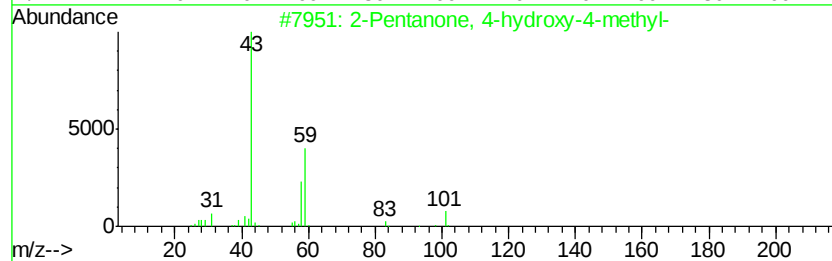
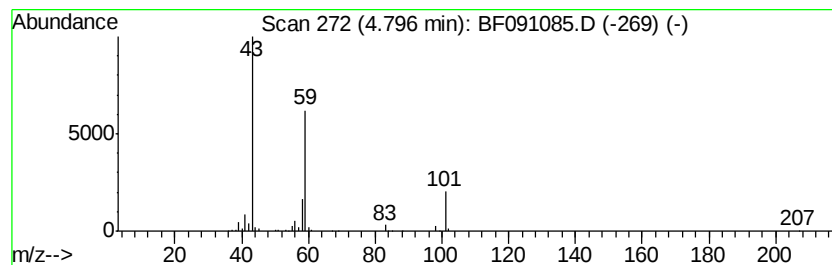
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.80	10.35 ng	576998	1,4-Dichlorobenzene-d4	6.66

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2		Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
3		Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	23
4		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
5		Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9



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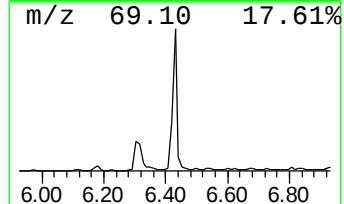
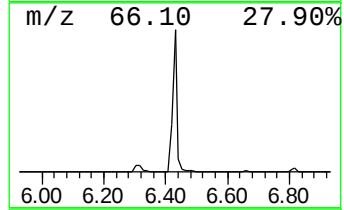
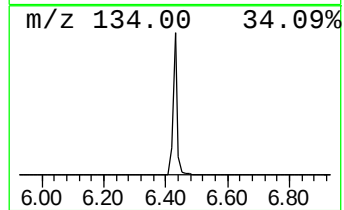
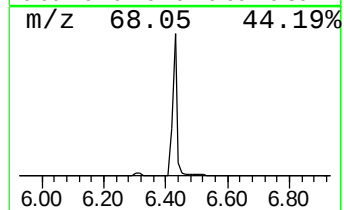
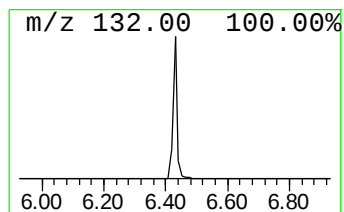
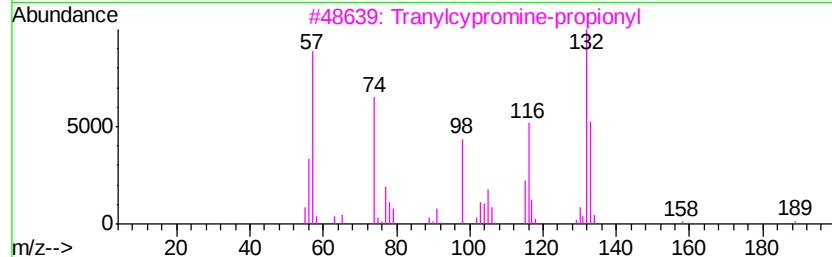
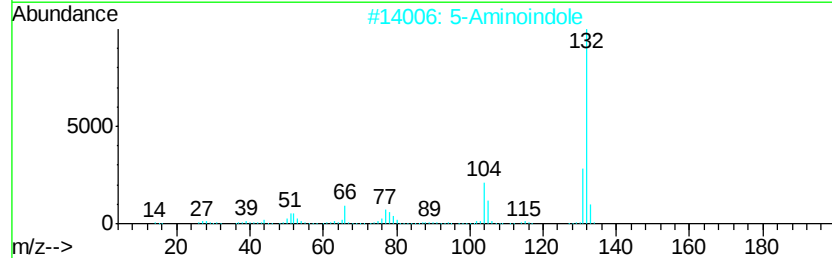
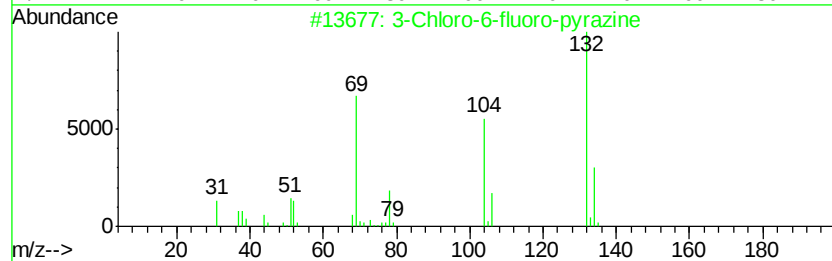
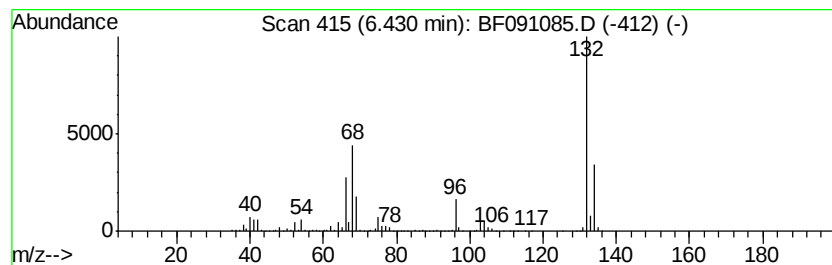
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 unknown6.43 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.43	73.37 ng	4088370	1,4-Dichlorobenzene-d4	6.66

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	3	Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16
2	5	Aminoindole	132	C8H8N2	005192-03-0	12
3		Tranvlcyproline-propionyl	189	C12H15NO	1000123-86-3	10
4		4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	9
5		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9



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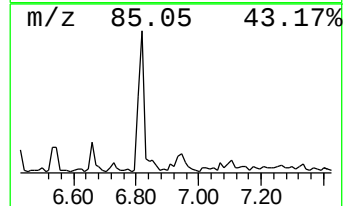
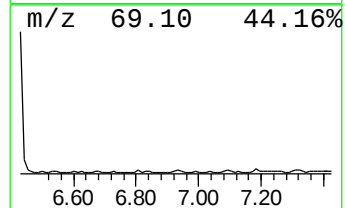
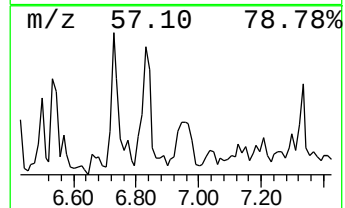
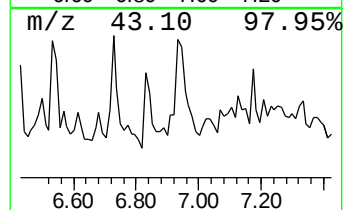
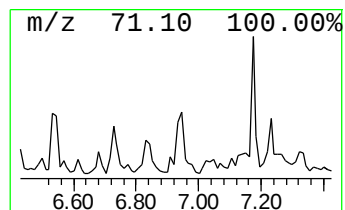
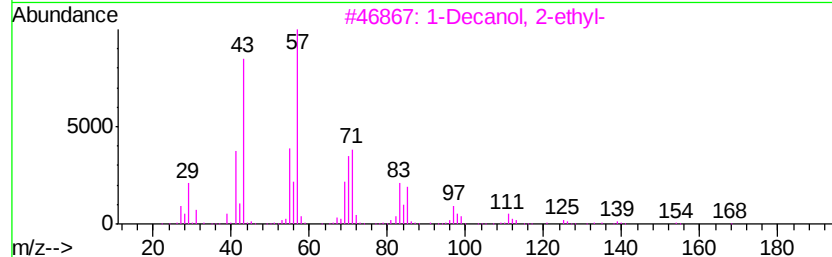
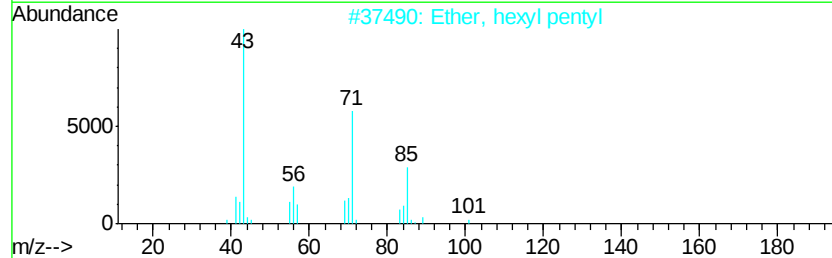
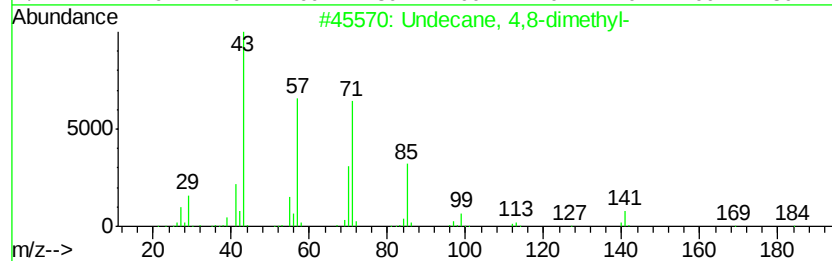
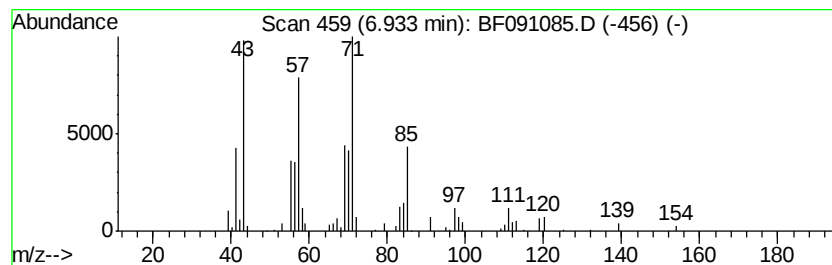
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 5 Undecane, 4,8-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.93	2.94 ng	163624	1,4-Dichlorobenzene-d4	6.66

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Undecane, 4,8-dimethyl-	184	C13H28	017301-33-6	64
2		Ether, hexyl pentyl	172	C11H24O	032357-83-8	59
3		1-Decanol, 2-ethyl-	186	C12H26O	021078-65-9	47
4		1-Decene, 4-methyl-	154	C11H22	013151-29-6	43
5		2,4,4-Trimethyl-1-hexene	126	C9H18	051174-12-0	43



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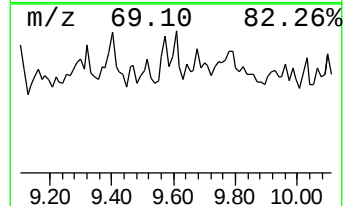
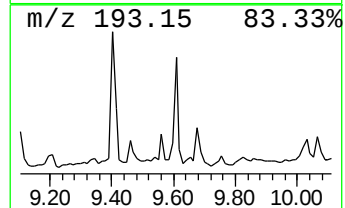
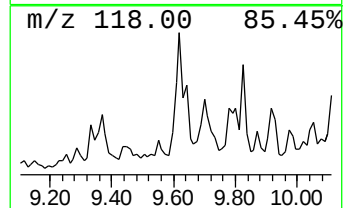
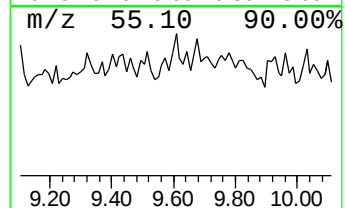
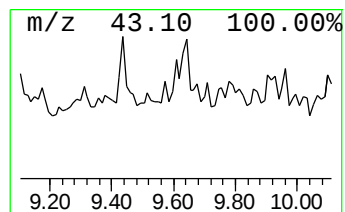
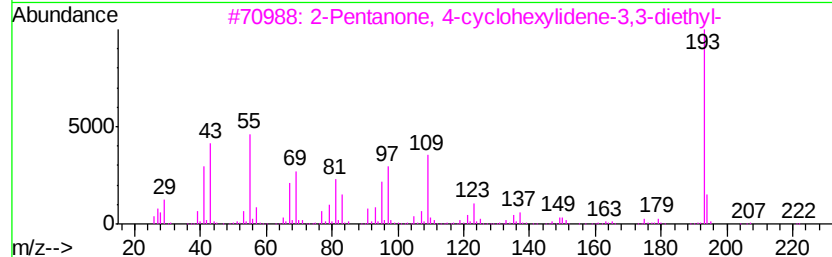
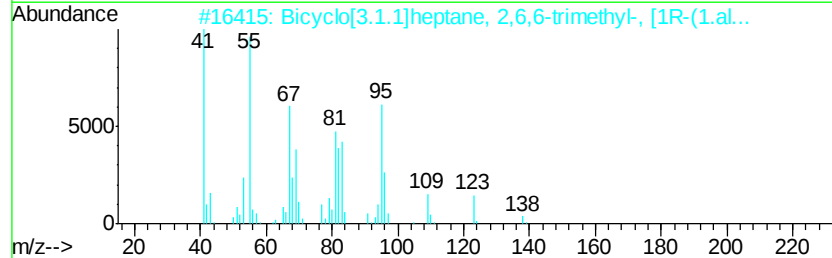
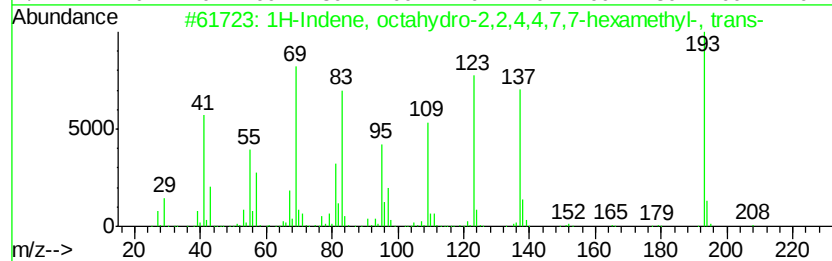
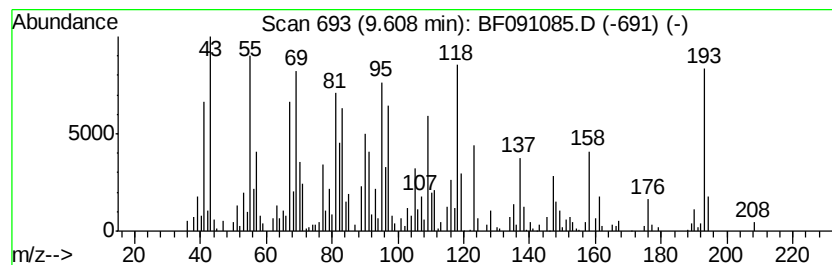
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 6 unknown9.61 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.61	2.12 ng	209030	Acenaphthene-d10	9.70

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Indene, octahydro-2,2,4,4,7,7...	208	C15H28	054832-83-6	35
2		Bicyclo[3.1.1]heptane, 2,6,6-tri...	138	C10H18	004863-59-6	25
3		2-Pentanone, 4-cyclohexylidene-3...	222	C15H26O	313253-65-5	22
4		2,7-Octadiene, 4-methyl-	124	C9H16	1000061-78-0	20
5		6-Acetamido-1,4-benzodioxane	193	C10H11NO3	063546-19-0	14



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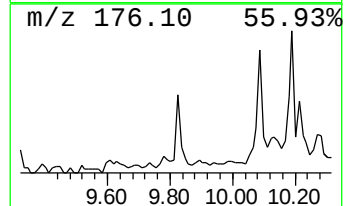
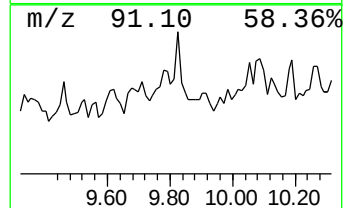
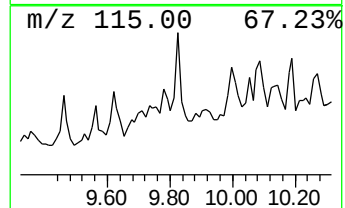
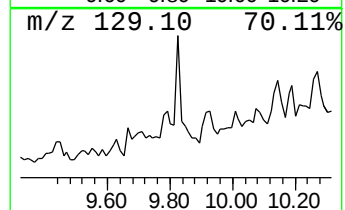
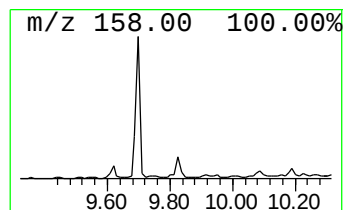
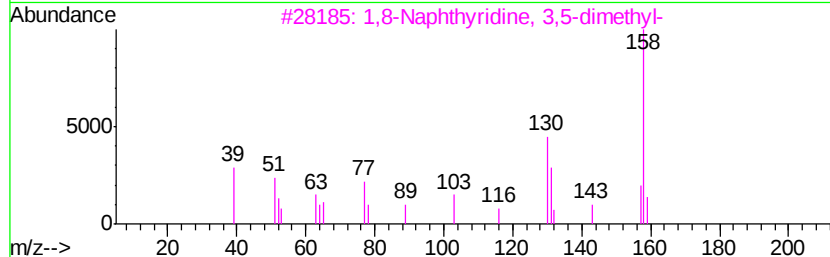
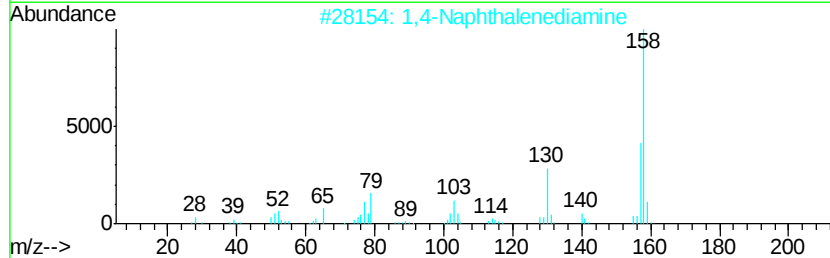
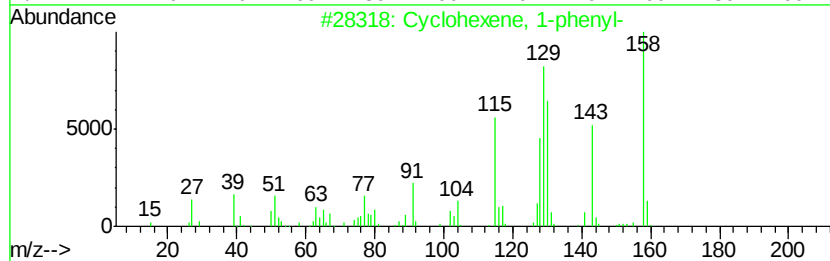
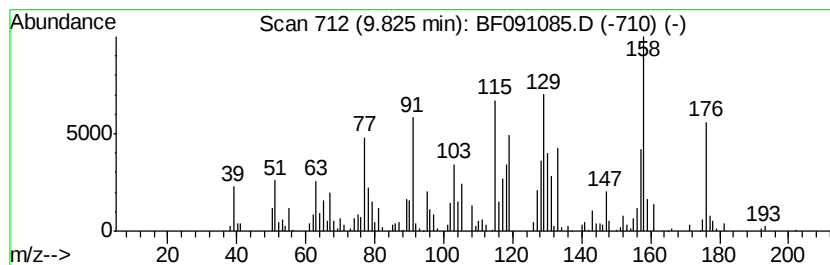
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Peak Number 7 unknown9.82 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.82	2.45 ng	241608	Acenaphthene-d10	9.70

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexene, 1-phenyl-	158	C12H14	000771-98-2	46
2		1,4-Naphthalenediamine	158	C10H10N2	002243-61-0	46
3		1,8-Naphthyridine, 3,5-dimethyl-	158	C10H10N2	014757-46-1	42
4		1H-Pyrazole, 3-methyl-5-phenyl-	158	C10H10N2	003347-62-4	41
5		7-Methyl-1-naphthol	158	C11H10O	006939-33-9	41



Data Path : Z:\HPCHEM1\BNA F\DATA\BF110816\
Data File : BF091085.D
Acq On : 8 Nov 2016 16:43
Operator : UM/SJ
Sample : H5571-01
Misc :
ALS Vial : 6 Sample Multiplier: 1

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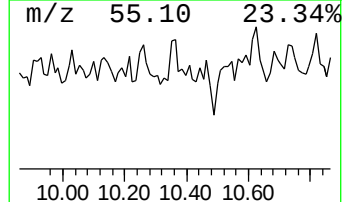
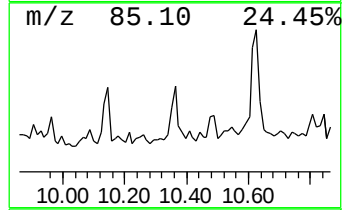
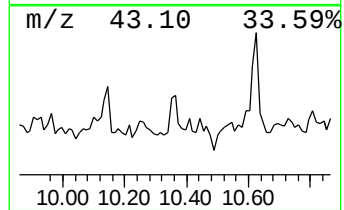
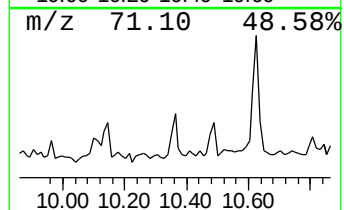
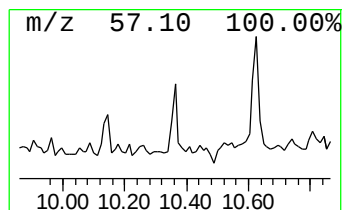
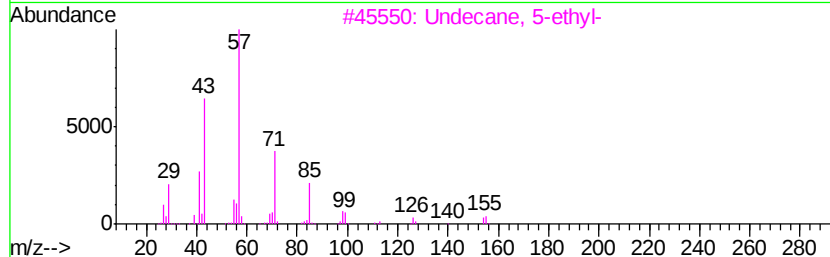
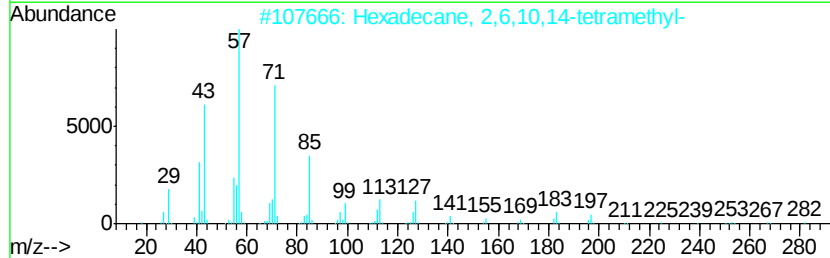
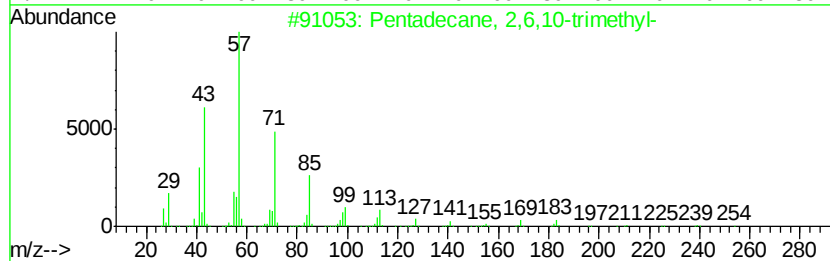
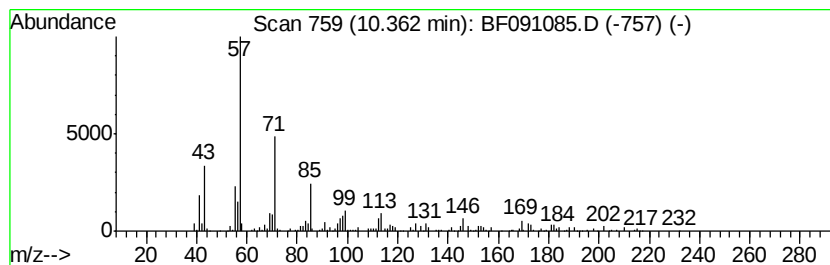
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Pentadecane, 2,6,10-trimethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.36	2.53 ng	249746	Acenaphthene-d10	9.70

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentadecane, 2,6,10-trimethyl-	254	C18H38	003892-00-0	72
2		Hexadecane, 2,6,10,14-tetramethyl-	282	C20H42	000638-36-8	72
3		Undecane, 5-ethyl-	184	C13H28	017453-94-0	64
4		Octadecane	254	C18H38	000593-45-3	64
5		Tridecane	184	C13H28	000629-50-5	64



Data Path : Z:\HPCHEM1\BNA F\DATA\BF110816\
Data File : BF091085.D
Acq On : 8 Nov 2016 16:43
Operator : UM/SJ
Sample : H5571-01
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
FORMER-POTABLE-WELL

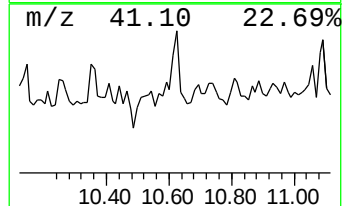
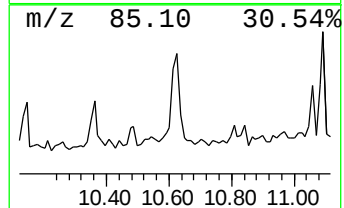
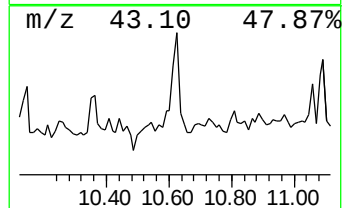
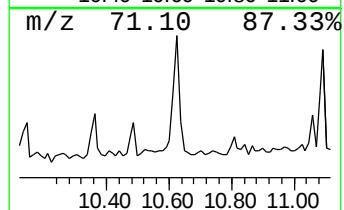
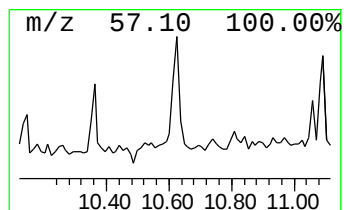
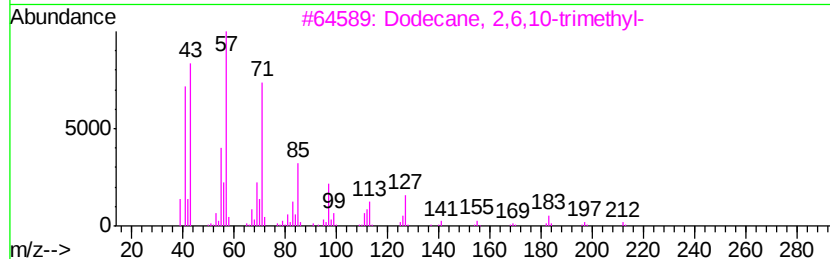
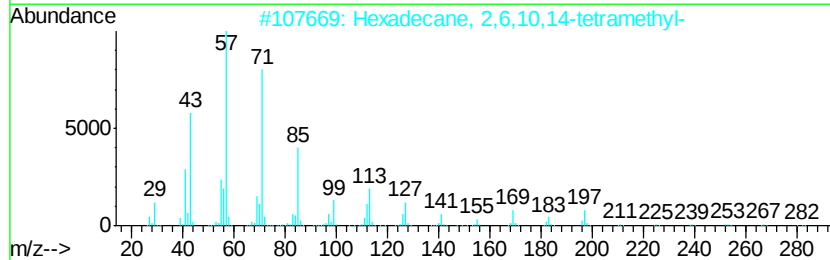
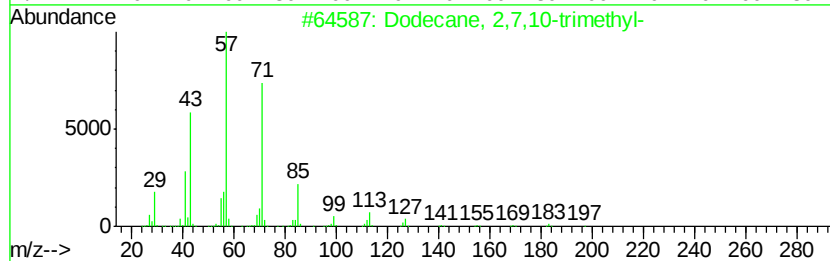
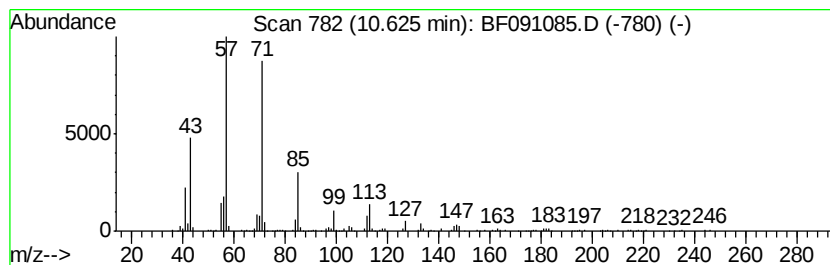
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Dodecane, 2,7,10-trimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.62	4.30 ng	334816	Phenanthrene-d10	11.17

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dodecane, 2,7,10-trimethyl-	212	C15H32	074645-98-0	86	
2	Hexadecane, 2,6,10,14-tetramethyl-	282	C20H42	000638-36-8	78	
3	Dodecane, 2,6,10-trimethyl-	212	C15H32	003891-98-3	78	
4	Undecane, 6-ethyl-	184	C13H28	017312-60-6	64	
5	Octane, 2,3,6-trimethyl-	156	C11H24	062016-33-5	59	



Data Path : Z:\HPCHEM1\BNA F\DATA\BF110816\
Data File : BF091085.D
Acq On : 8 Nov 2016 16:43
Operator : UM/SJ
Sample : H5571-01
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
FORMER-POTABLE-WELL

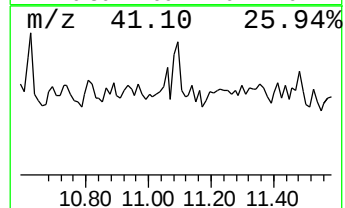
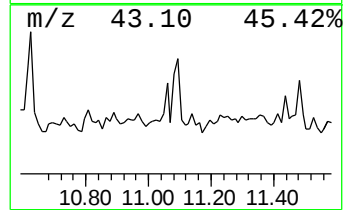
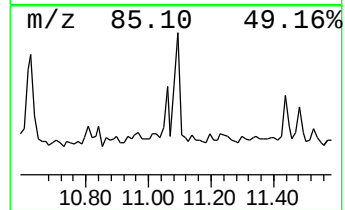
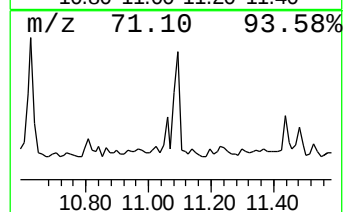
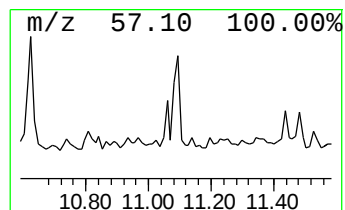
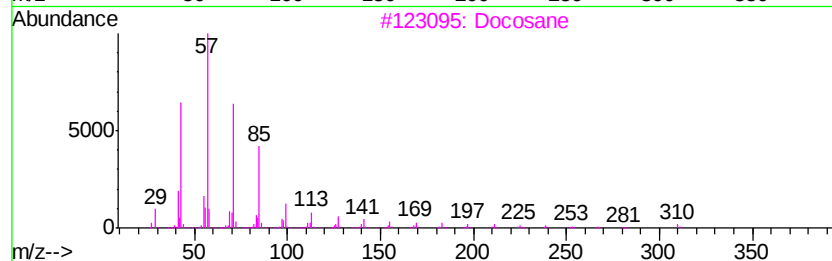
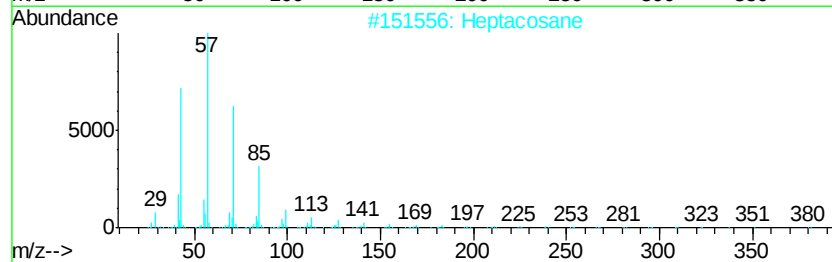
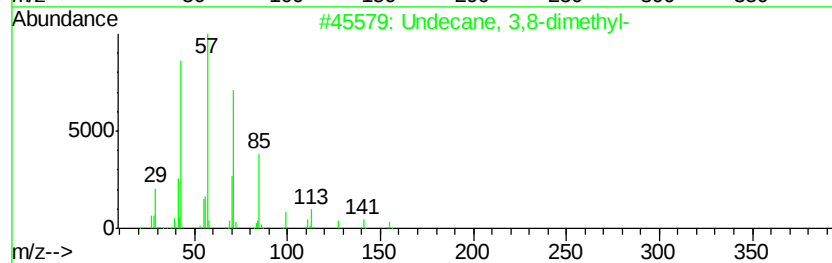
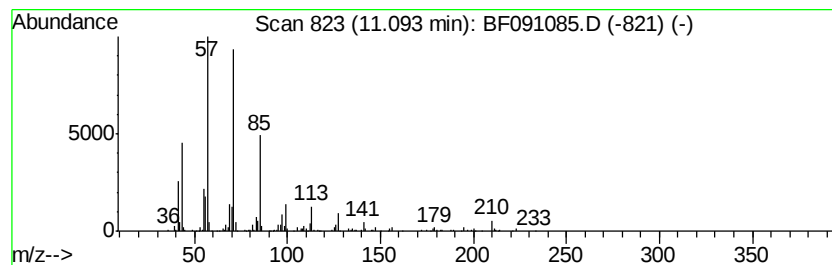
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Undecane, 3,8-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.09	3.15 ng	245201	Phenanthrene-d10	11.17

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Undecane, 3,8-dimethyl-	184	C13H28	017301-30-3	83
2		Heptacosane	380	C27H56	000593-49-7	80
3		Docosane	310	C22H46	000629-97-0	78
4		Dodecane, 2,6,11-trimethyl-	212	C15H32	031295-56-4	72
5		Heneicosane	296	C21H44	000629-94-7	64



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF110816\
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Acq On : 8 Nov 2016 16:43
Operator : UM/SJ
Sample : H5571-01
Misc :
ALS Vial : 6 Sample Multiplier: 1

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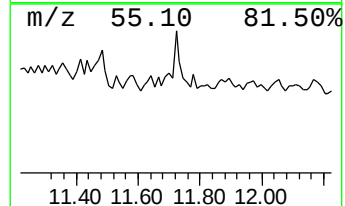
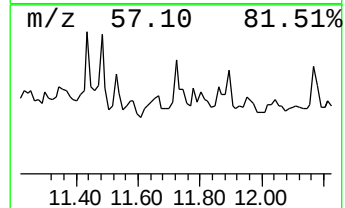
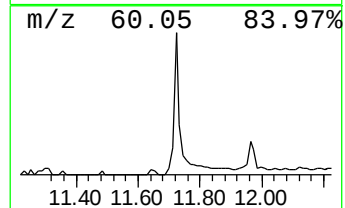
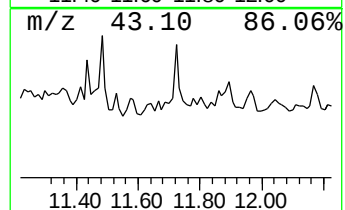
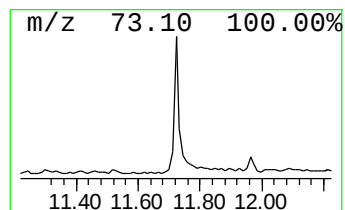
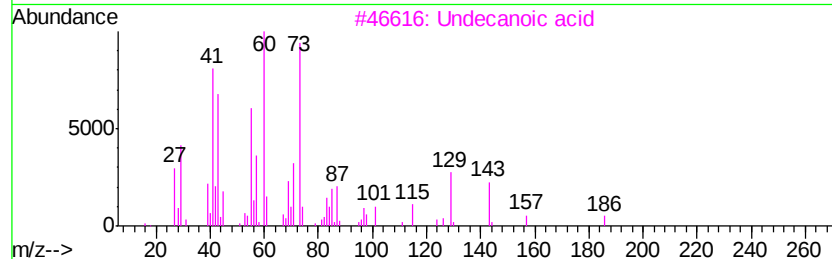
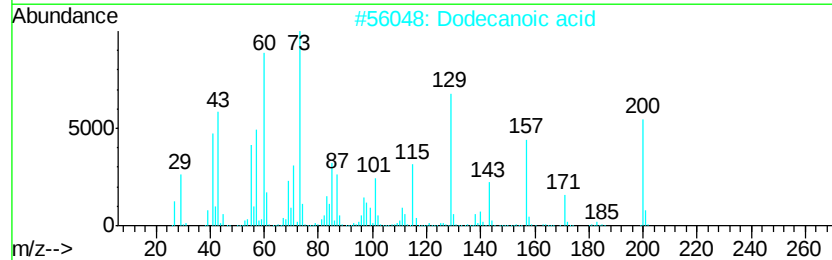
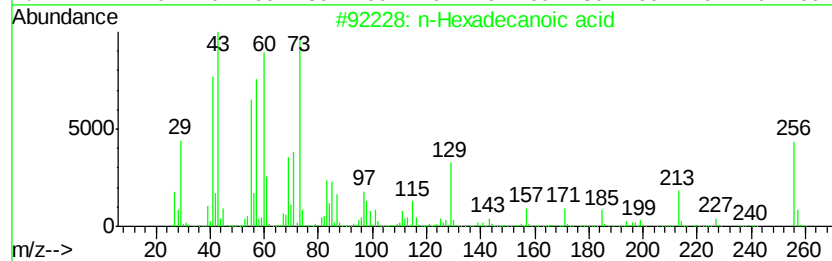
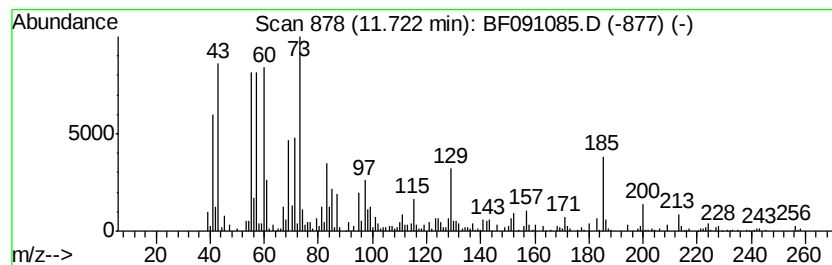
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 n-Hexadecanoic acid Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.72	2.89 ng	224966	Phenanthrene-d10	11.17

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95
2		Dodecanoic acid	200	C12H24O2	000143-07-7	55
3		Undecanoic acid	186	C11H22O2	000112-37-8	52
4		Tridecanoic acid	214	C13H26O2	000638-53-9	50
5		Pentadecanoic acid	242	C15H30O2	001002-84-2	38



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF110816\
Data File : BF091085.D
Acq On : 8 Nov 2016 16:43
Operator : UM/SJ
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Misc :
ALS Vial : 6 Sample Multiplier: 1

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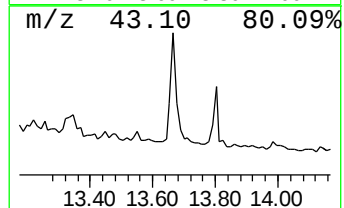
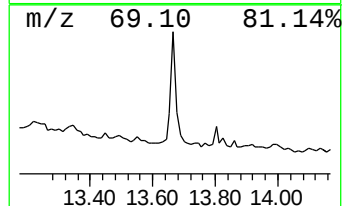
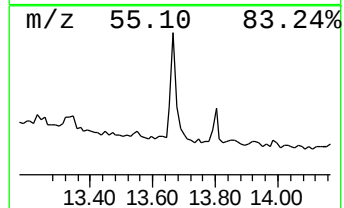
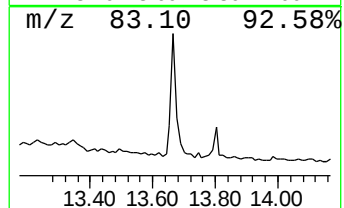
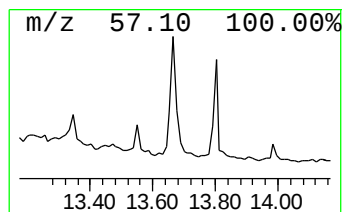
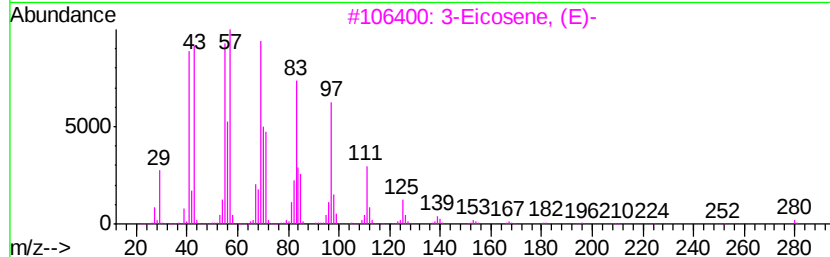
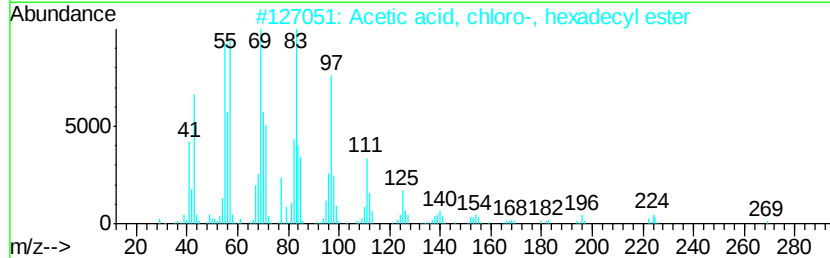
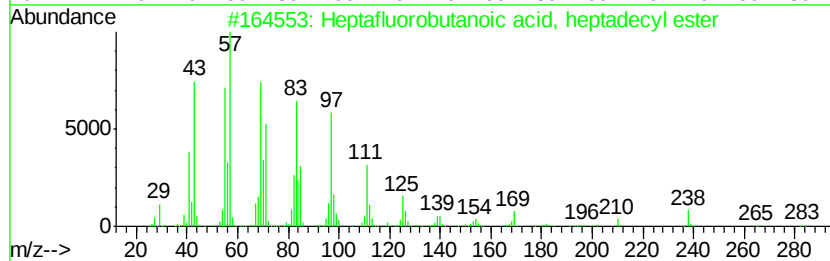
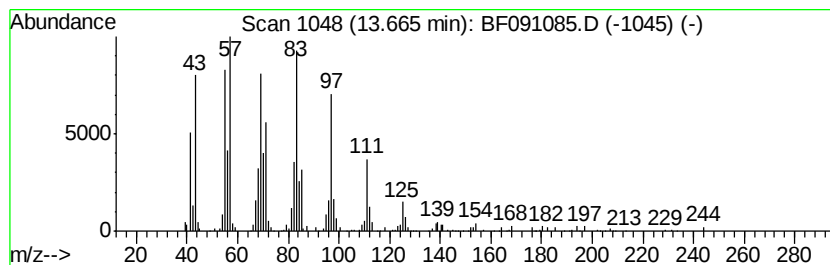
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Heptafluorobutanoic acid, h... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.67	2.93 ng	246815	Chrysene-d12	13.80

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptafluorobutanoic acid, heptad...	452	C21H35F7O2	1000282-97-3	93
2		Acetic acid, chloro-, hexadecyl ...	318	C18H35ClO2	052132-58-8	91
3		3-Eicosene, (E)-	280	C20H40	074685-33-9	91
4		1-Nonadecene	266	C19H38	018435-45-5	91
5		Dichloroacetic acid, heptadecyl ...	366	C19H36Cl2O2	1000282-98-2	91



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF110816\
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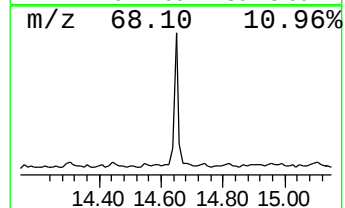
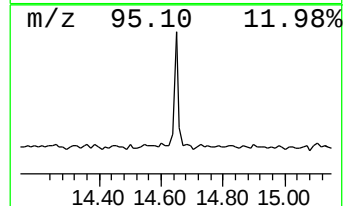
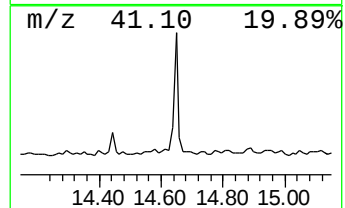
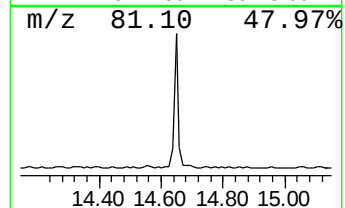
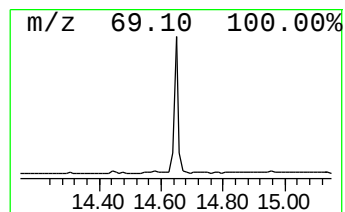
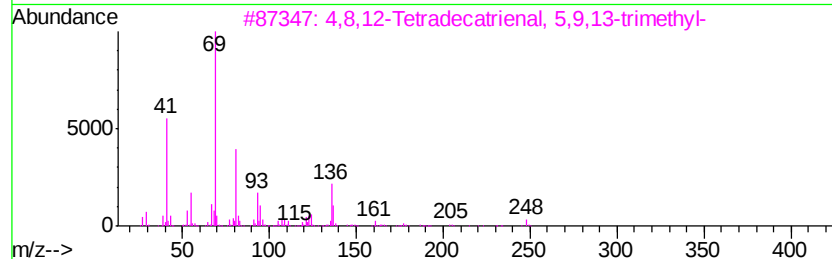
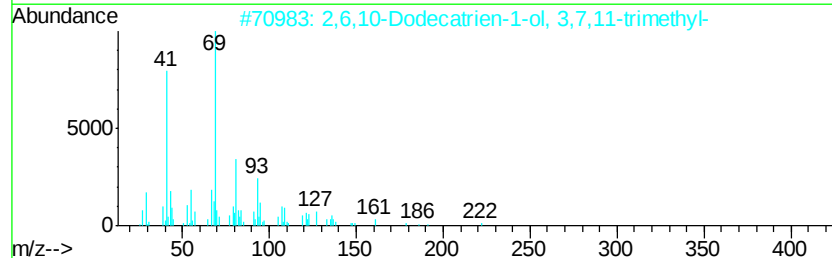
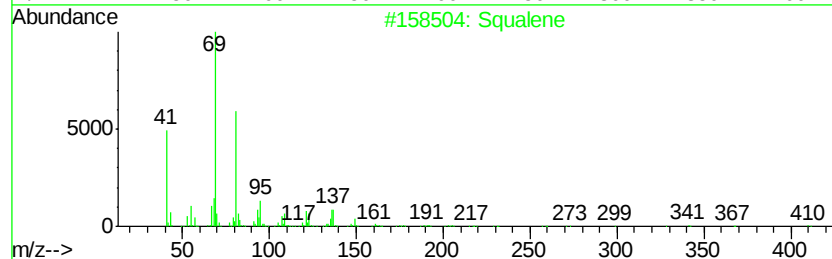
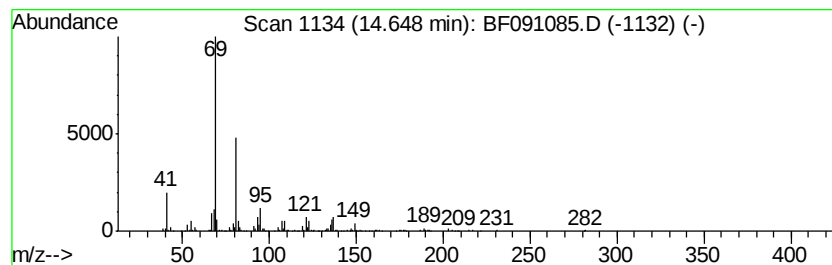
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Squalene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.65	3.62 ng	210481	Perylene-d12	15.19

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Squalene	410	C30H50	007683-64-9	91
2		2,6,10-Dodecatrien-1-ol, 3,7,11-...	222	C15H26O	004602-84-0	72
3		4,8,12-Tetradecatrienal, 5,9,13-...	248	C17H28O	066408-55-7	72
4		7,11-Dimethyldodeca-2,6,10-trien...	208	C14H24O	125529-10-4	64
5		2,3,5-Trimethyl-2,3,5-hexanetric...	203	C12H17N3	003974-79-6	64



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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF110316.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
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unknown6.43	6.43	73.4	ng	4088370	1	6.66	1114440	20.0
Undecane, 4,8-dim...	6.93	2.9	ng	163624	1	6.66	1114440	20.0
unknown9.61	9.61	2.1	ng	209030	3	9.70	1972870	20.0
unknown9.82	9.82	2.5	ng	241608	3	9.70	1972870	20.0
Pentadecane, 2,6,...	10.36	2.5	ng	249746	3	9.70	1972870	20.0
Dodecane, 2,7,10-...	10.62	4.3	ng	334816	4	11.17	1558530	20.0
Undecane, 3,8-dim...	11.09	3.1	ng	245201	4	11.17	1558530	20.0
n-Hexadecanoic acid	11.72	2.9	ng	224966	4	11.17	1558530	20.0
Heptafluorobutano...	13.67	2.9	ng	246815	5	13.80	1681910	20.0
Squalene	14.65	3.6	ng	210481	6	15.19	1162670	20.0