

Data Path : Z:\HPCHEM1\BNA F\DATA\BF042016\
 Data File : BF086318.D
 Acq On : 20 Apr 2016 19:43
 Operator : UM/SJ
 Sample : H2625-05
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW-2

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-BF042016.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	2.178	5	8	41	rVB5	136272	914282	13.43%	1.730%
2	3.058	81	85	95	rVB	68502	154603	2.27%	0.292%
3	3.607	124	133	138	rVB	41619	111178	1.63%	0.210%
4	5.047	256	259	261	rBV	1149314	1331457	19.55%	2.519%
5	5.459	292	295	297	rBV	2348382	2373361	34.85%	4.490%
6	5.870	328	331	333	rBV	208698	245782	3.61%	0.465%
7	6.476	382	384	394	rBV	1031034	1805579	26.52%	3.416%
8	6.624	394	397	410	rVV	4235936	4573611	67.16%	8.652%
9	6.864	415	418	420	rBV	951862	1367925	20.09%	2.588%
10	7.013	429	431	434	rBV	3789325	4348419	63.86%	8.226%
11	7.185	444	446	451	rVB4	53868	90565	1.33%	0.171%
12	7.425	460	467	473	rBV	3664234	5169678	75.92%	9.780%
13	7.505	473	474	479	rVV4	94940	233390	3.43%	0.442%
14	7.630	483	485	489	rVV3	79626	134393	1.97%	0.254%
15	7.710	489	492	494	rVV2	98118	171639	2.52%	0.325%
16	7.836	497	503	504	rBV6	28870	69173	1.02%	0.131%
17	7.882	504	507	510	rBV5	35482	68724	1.01%	0.130%
18	8.042	517	521	524	rBV2	90824	142154	2.09%	0.269%
19	8.145	527	530	537	rBV	1828383	2045518	30.04%	3.870%
20	8.693	574	578	581	rBV4	38812	85511	1.26%	0.162%
21	8.785	583	586	587	rBV	46980	76432	1.12%	0.145%
22	9.230	621	625	627	rBV	4715453	6809521	100.00%	12.882%
23	9.619	657	659	661	rBV	260664	271681	3.99%	0.514%
24	9.836	676	678	681	rBV	126553	122167	1.79%	0.231%
25	9.893	681	683	686	rVV	1365902	2031266	29.83%	3.843%
26	10.271	708	716	719	rBV	155513	750529	11.02%	1.420%
27	10.339	719	722	723	rVV2	255447	388348	5.70%	0.735%
28	10.362	723	724	735	rVB7	58483	192742	2.83%	0.365%
29	10.556	738	741	744	rBV	84111	140671	2.07%	0.266%
30	10.682	750	752	761	rBV	2144732	2894953	42.51%	5.476%
31	10.808	761	763	770	rVB	330743	432438	6.35%	0.818%
32	11.254	799	802	804	rBV2	139099	160806	2.36%	0.304%
33	11.379	811	813	817	rBV	1683344	1841257	27.04%	3.483%
34	11.562	827	829	831	rBV	45126	76709	1.13%	0.145%

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

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
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Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	11.688	837	840	846	rVB2	52630	135158	1.98%	0.256%
36	12.797	934	937	940	rBV2	44227	69609	1.02%	0.132%
37	12.865	940	943	946	rBV2	217638	333987	4.90%	0.632%
38	12.968	949	952	958	rBV	4194294	4799685	70.48%	9.080%
39	13.208	970	973	974	rBV	90346	131821	1.94%	0.249%
40	13.539	1000	1002	1007	rVB4	214380	373123	5.48%	0.706%
41	13.871	1029	1031	1033	rBV	378371	354051	5.20%	0.670%
42	13.917	1033	1035	1041	rVV	238588	709627	10.42%	1.342%
43	14.008	1041	1043	1051	rVB	1064362	1420727	20.86%	2.688%
44	14.191	1056	1059	1061	rBV	395459	383394	5.63%	0.725%
45	14.488	1083	1085	1088	rBV	267817	313494	4.60%	0.593%
46	14.785	1109	1111	1114	rBV	140983	209603	3.08%	0.397%
47	15.117	1137	1140	1145	rBV	107906	126520	1.86%	0.239%
48	15.437	1165	1168	1171	rBV	844139	1190673	17.49%	2.252%
49	15.482	1171	1172	1185	rVB2	69007	111589	1.64%	0.211%
50	16.054	1217	1222	1226	rBV8	19338	83357	1.22%	0.158%
51	16.385	1247	1251	1260	rVB7	35686	109633	1.61%	0.207%
52	17.494	1342	1348	1352	rBV7	59819	258700	3.80%	0.489%
53	17.894	1380	1383	1391	rVB4	44373	121040	1.78%	0.229%

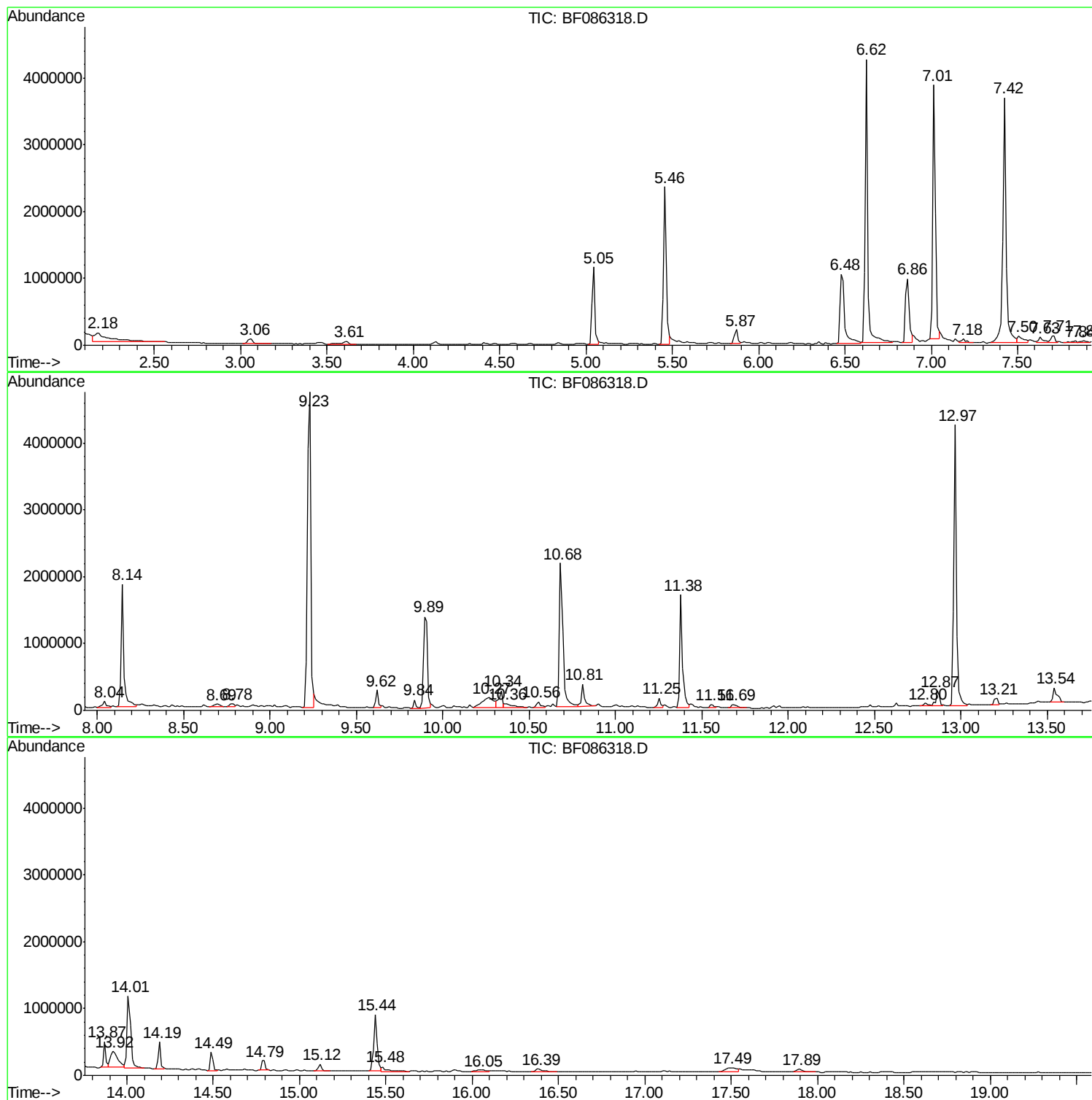
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BNA_F
ClientSampled :
MW-2

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF042016.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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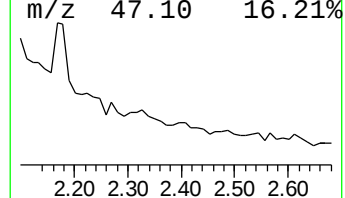
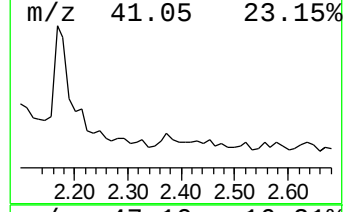
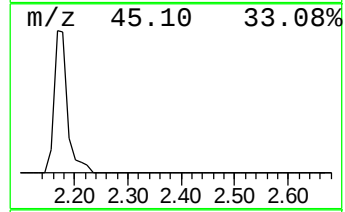
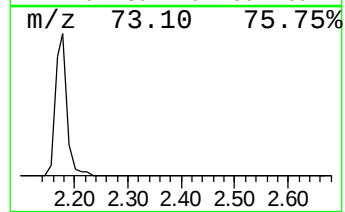
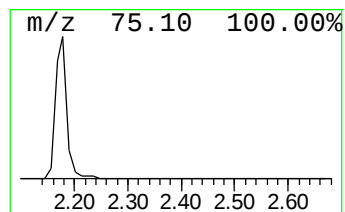
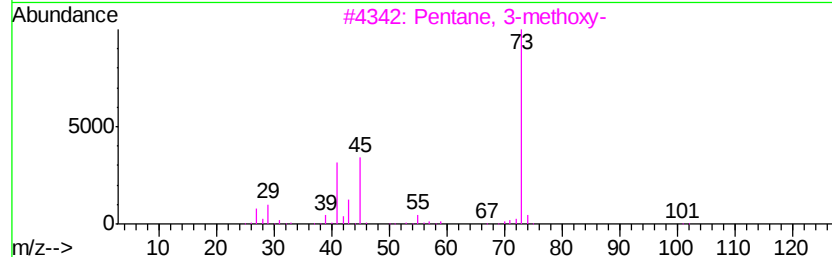
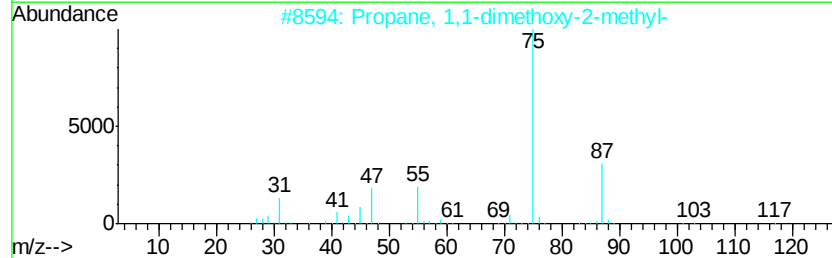
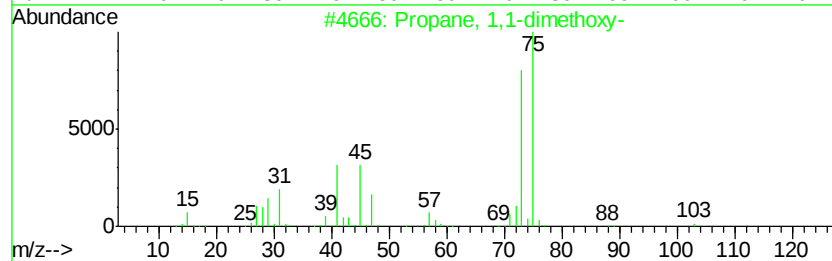
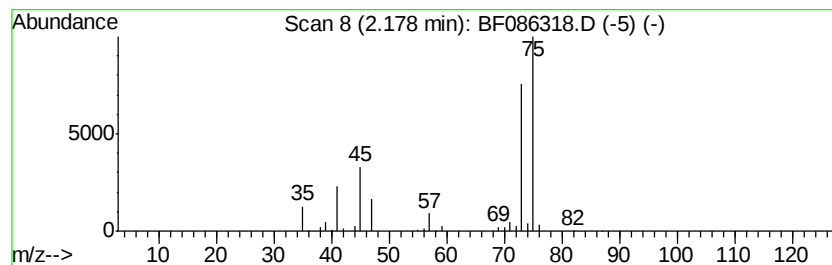
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF042016.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Propane, 1,1-dimethoxy- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.18	13.37 ng	914282	1,4-Dichlorobenzene-d4	6.86

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propane, 1,1-dimethoxy-	104	C5H12O2	004744-10-9	78
2			Propane, 1,1-dimethoxy-2-methyl-	118	C6H14O2	041632-89-7	28
3			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	14
4			Thiocyanic acid, methyl ester	73	C2H3NS	000556-64-9	12
5			Glycine	75	C2H5NO2	000056-40-6	9



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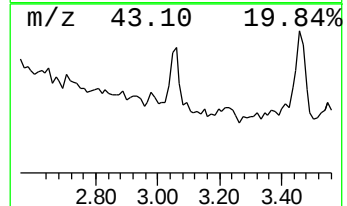
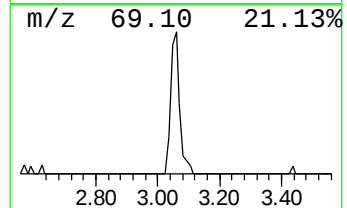
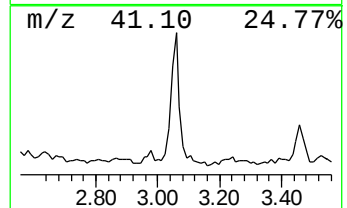
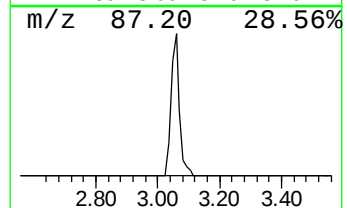
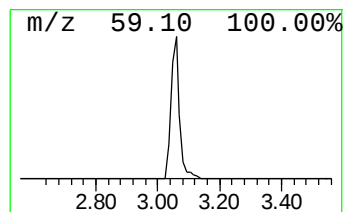
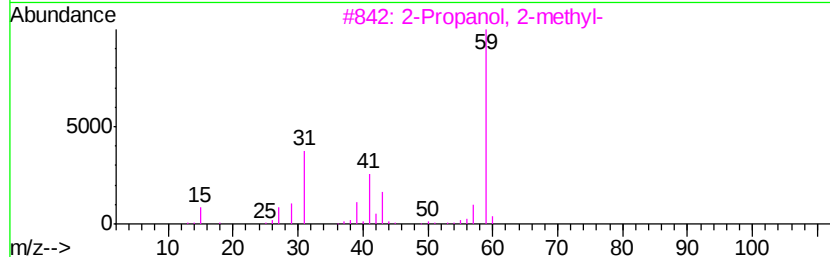
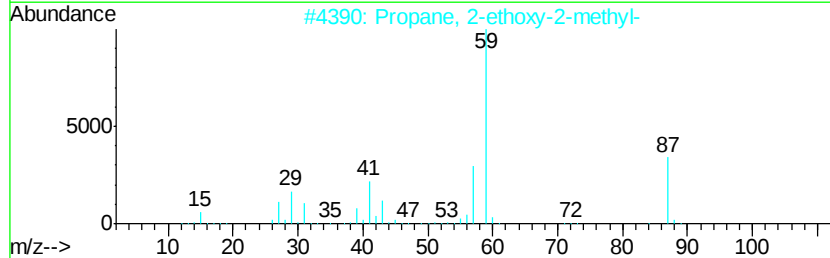
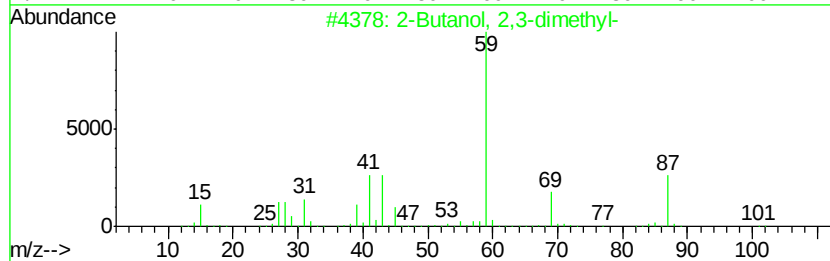
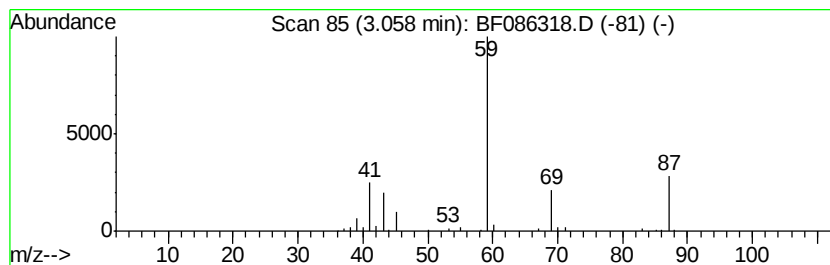
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 2-Butanol, 2,3-dimethyl- Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.06	2.26 ng	154603	1,4-Dichlorobenzene-d4	6.86

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Butanol, 2,3-dimethyl-	102	C6H14O	000594-60-5	78
2		Propane, 2-ethoxy-2-methyl-	102	C6H14O	000637-92-3	39
3		2-Propanol, 2-methyl-	74	C4H10O	000075-65-0	38
4		Amylene Hydrate	88	C5H12O	000075-85-4	36
5		N-Formyl-d-threo-0-methylthreonine	161	C6H11NO4	1000214-69-5	9



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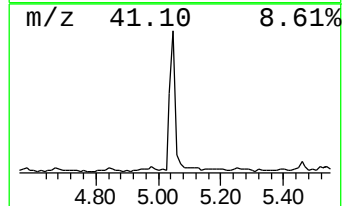
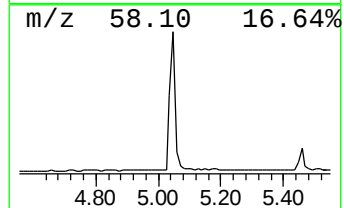
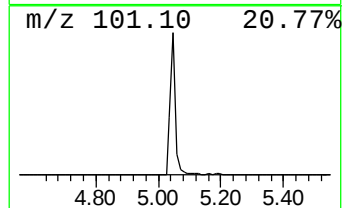
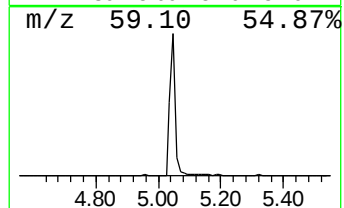
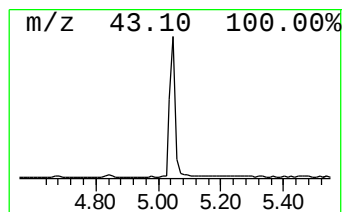
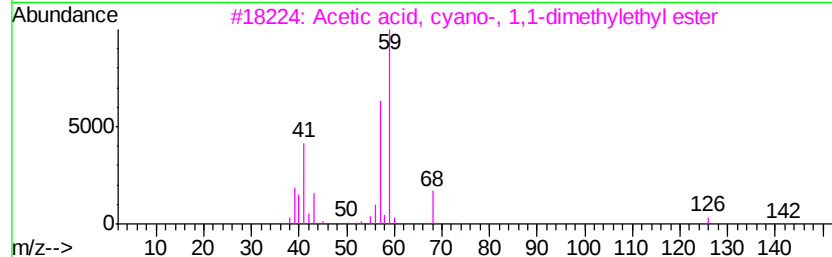
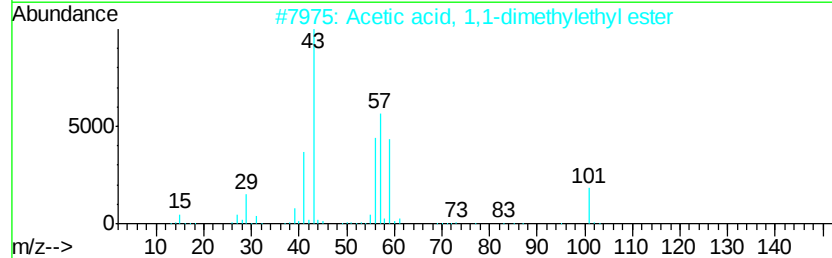
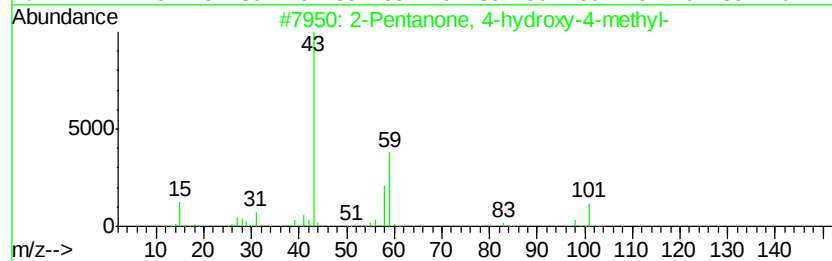
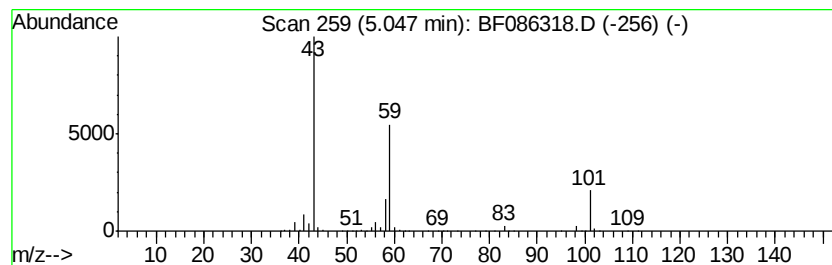
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.05	19.47 ng	1331460	1,4-Dichlorobenzene-d4	6.86

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	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		Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	17
4		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
5		Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9



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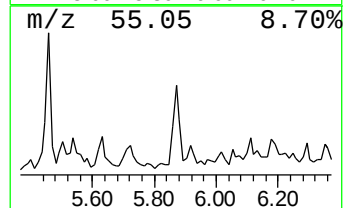
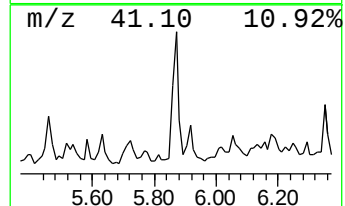
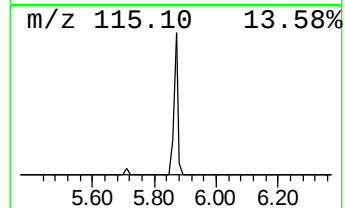
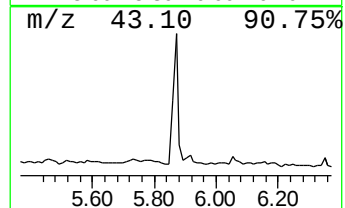
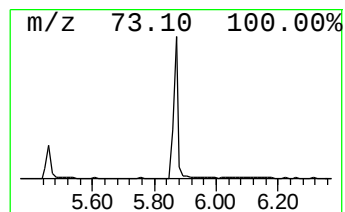
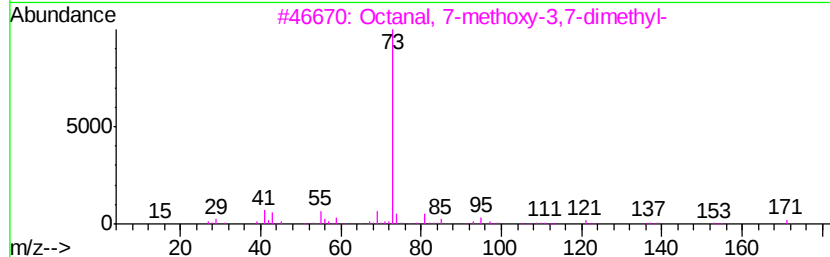
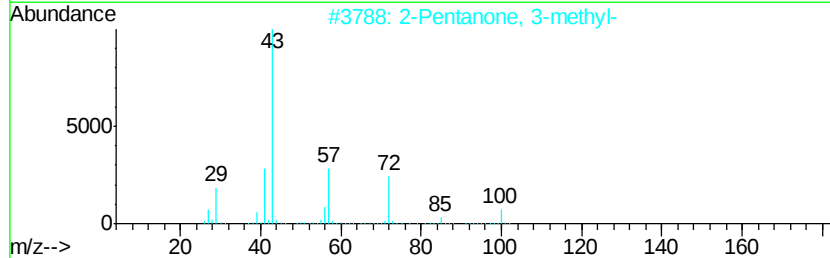
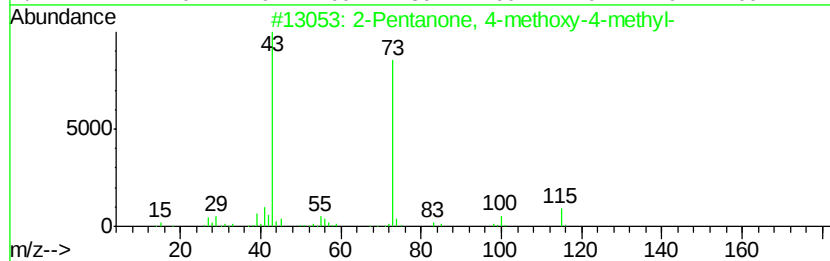
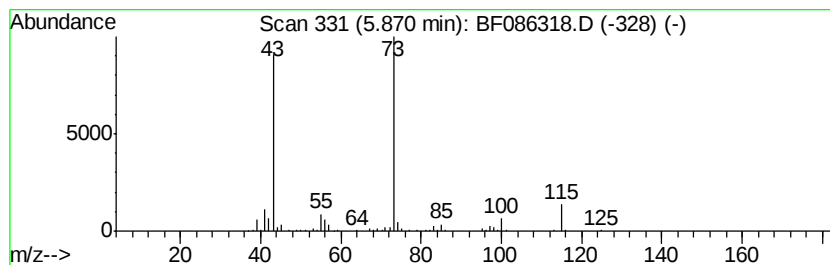
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 4 2-Pentanone, 4-methoxy-4-me... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.87	3.59 ng	245782	1,4-Dichlorobenzene-d4	6.86

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-methoxy-4-methyl-	130	C7H14O2	000107-70-0	78	
2	2-Pentanone, 3-methyl-	100	C6H12O	000565-61-7	27	
3	Octanal, 7-methoxy-3,7-dimethyl-	186	C11H22O2	003613-30-7	17	
4	1,3-Dioxolane, 2-pentadecyl-	284	C18H36O2	004360-57-0	16	
5	Heptane	100	C7H16	000142-82-5	12	



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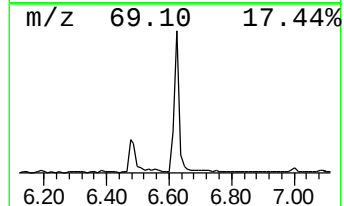
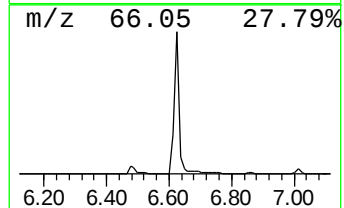
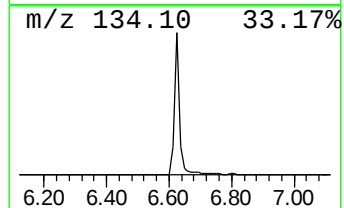
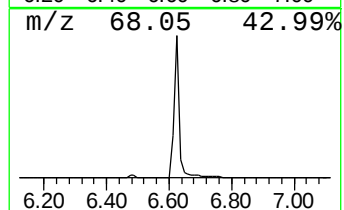
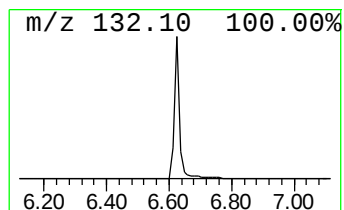
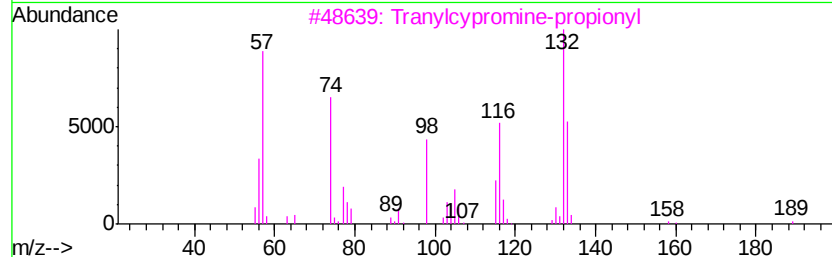
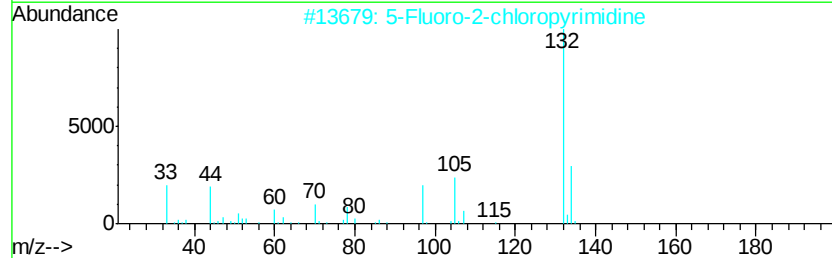
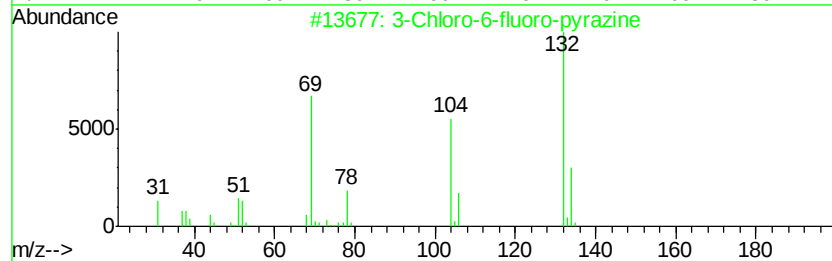
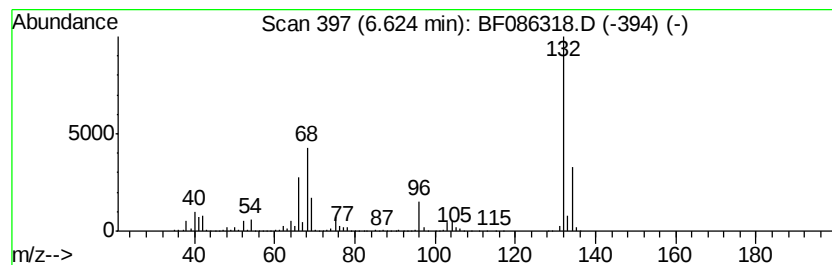
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 unknown6.62 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.62	66.87 ng	4573610	1,4-Dichlorobenzene-d4	6.86

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27	
2	5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16	
3	Tranvlcyproline-propionyl	189	C12H15NO	1000123-86-3	10	
4	Benzene, 1,2,4-trifluoro-	132	C6H3F3	000367-23-7	9	
5	Benzene, 1,3,5-trifluoro-	132	C6H3F3	000372-38-3	9	



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF042016\
Data File : BF086318.D
Acq On : 20 Apr 2016 19:43
Operator : UM/SJ
Sample : H2625-05
Misc :
ALS Vial : 22 Sample Multiplier: 1

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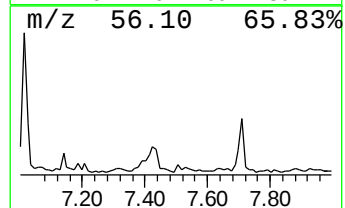
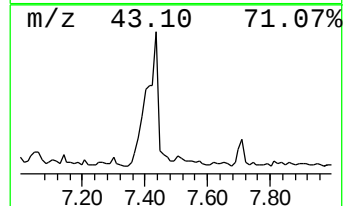
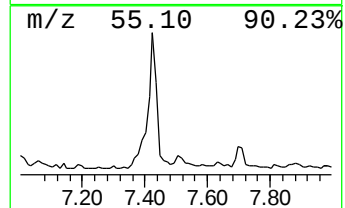
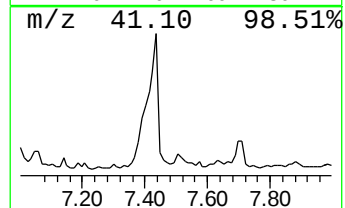
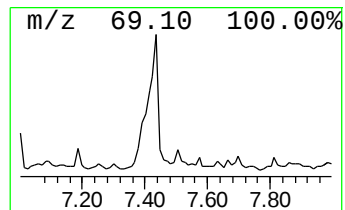
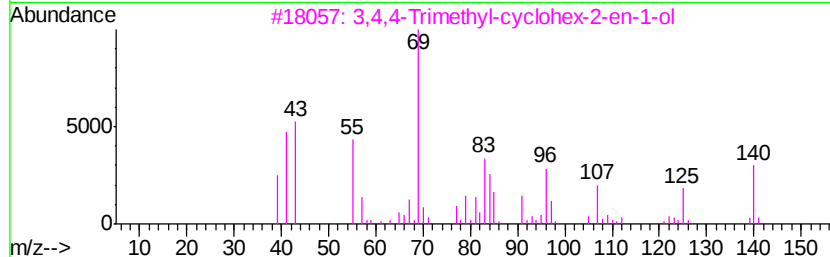
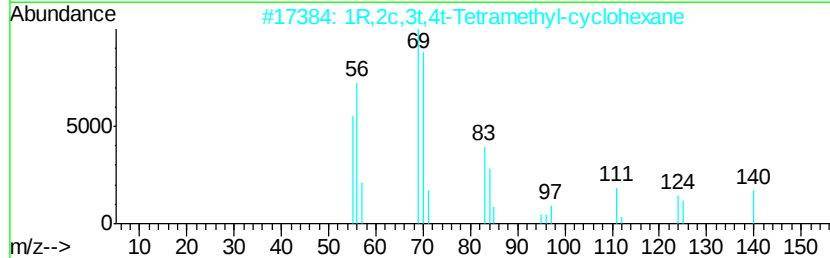
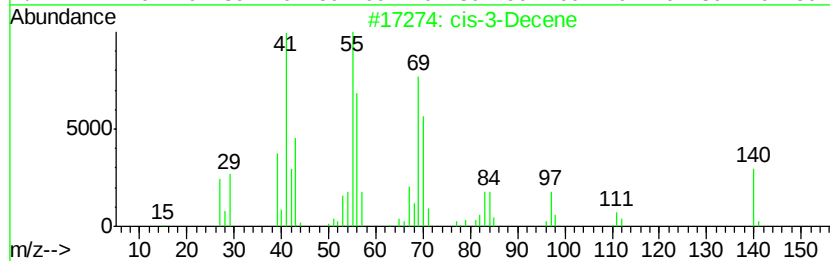
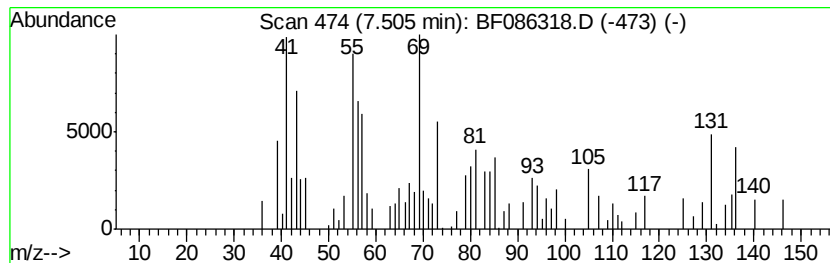
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 unknown7.50 Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.50	2.28 ng	233390	Naphthalene-d8	8.14

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	cis-3-Decene	140	C10H20	019398-86-8	46
2		1R,2c,3t,4t-Tetramethyl-cyclohexane	140	C10H20	1000144-07-3	46
3		3,4,4-Trimethyl-cyclohex-2-en-1-ol	140	C9H16O	073741-63-6	38
4		Cyclopropane, 1-ethyl-1-methyl-	84	C6H12	053778-43-1	38
5		1-Pentene, 3-methyl-	84	C6H12	000760-20-3	38



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Sample : H2625-05
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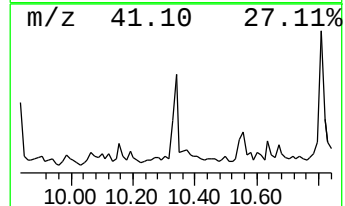
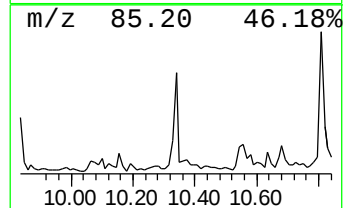
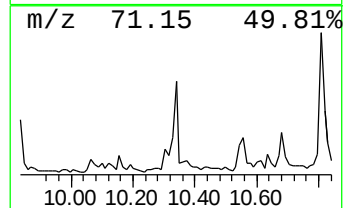
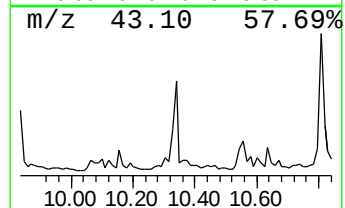
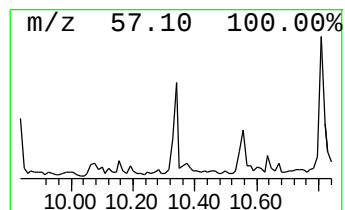
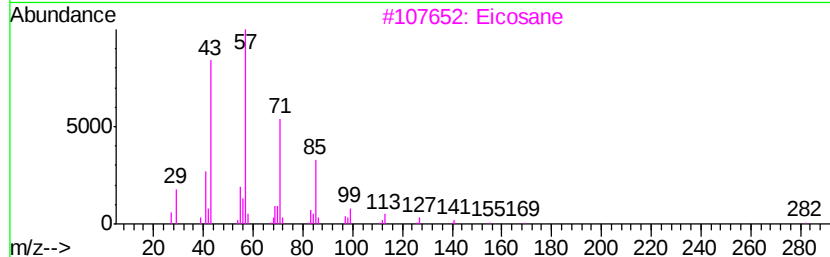
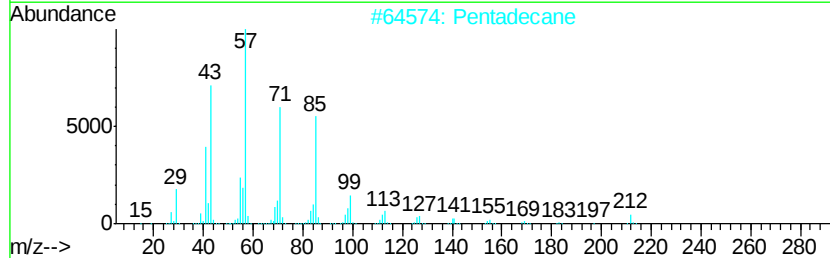
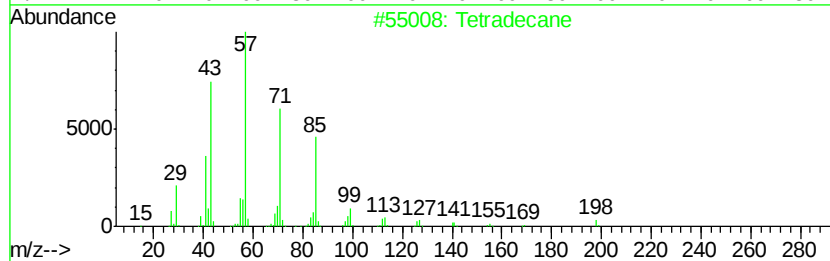
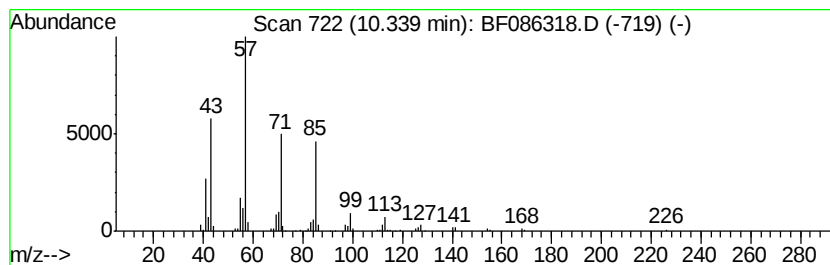
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 9 Tetradecane Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD		R.T.
10.34	3.82 ng	388348	Acenaphthene-d10		9.89
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetradecane	198	C14H30	000629-59-4	86
2	Pentadecane	212	C15H32	000629-62-9	86
3	Eicosane	282	C20H42	000112-95-8	83
4	Hexadecane	226	C16H34	000544-76-3	76
5	Decane, 3-bromo-	220	C10H21Br	030571-71-2	64



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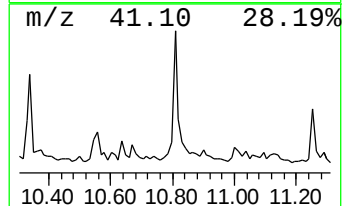
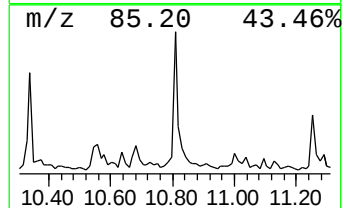
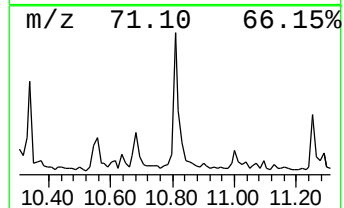
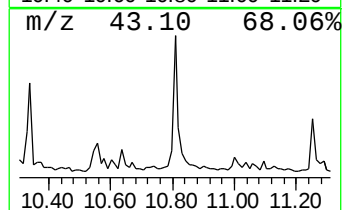
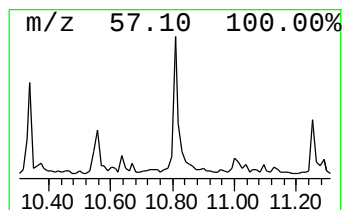
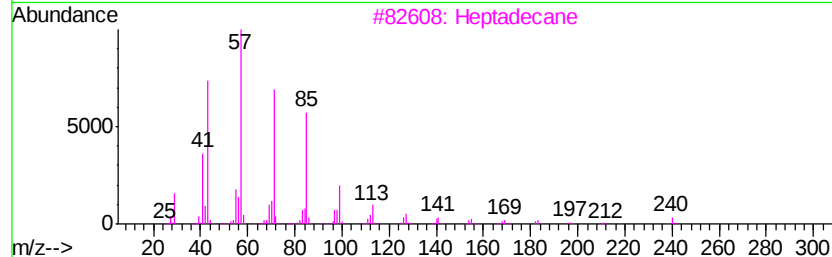
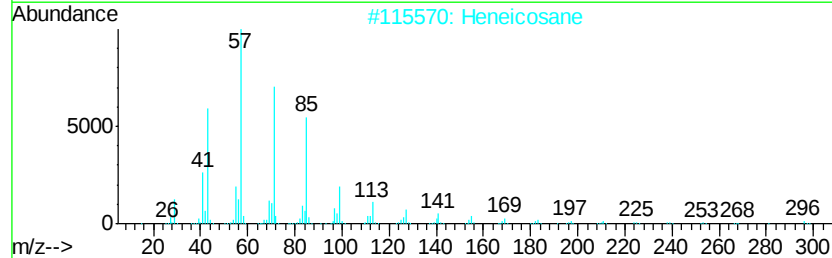
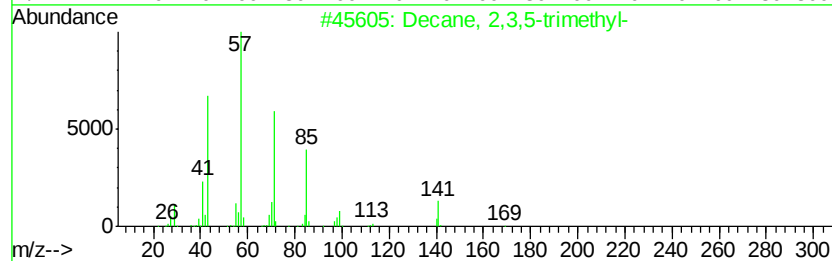
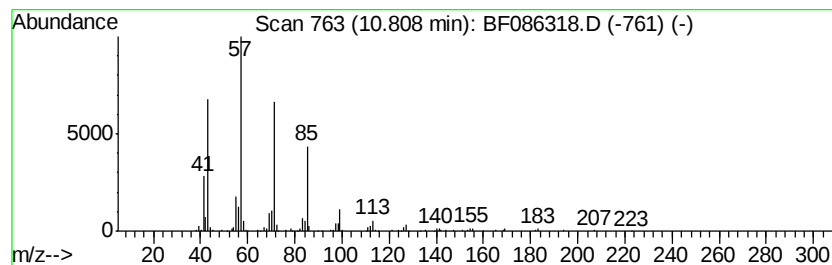
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 Decane, 2,3,5-trimethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.81	4.70 ng	432438	Phenanthrene-d10	11.38

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Decane, 2,3,5-trimethyl-	184	C13H28	062238-11-3	90
2			Heneicosane	296	C21H44	000629-94-7	90
3			Heptadecane	240	C17H36	000629-78-7	90
4			Eicosane	282	C20H42	000112-95-8	83
5			Hexadecane	226	C16H34	000544-76-3	72



Data Path : Z:\HPCHEM1\BNA F\DATA\BF042016\
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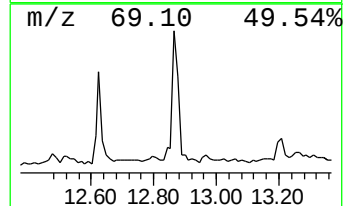
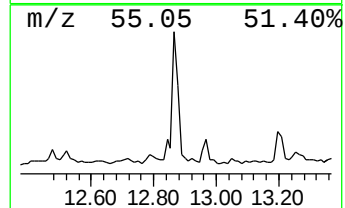
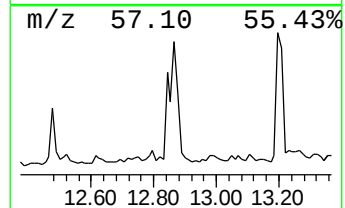
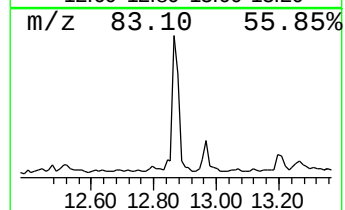
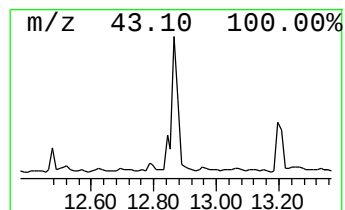
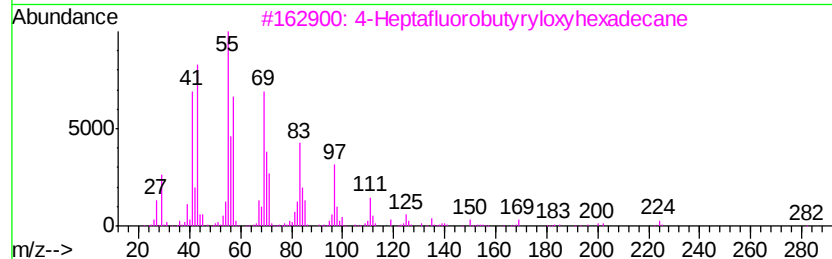
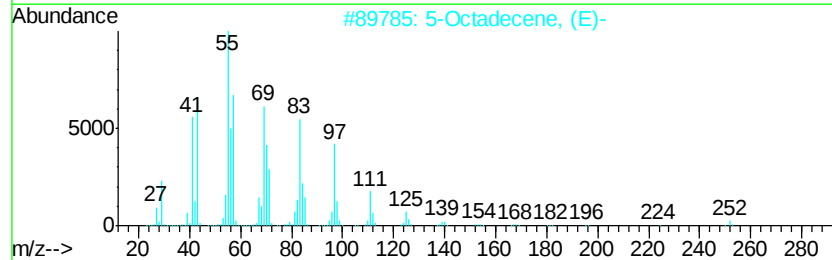
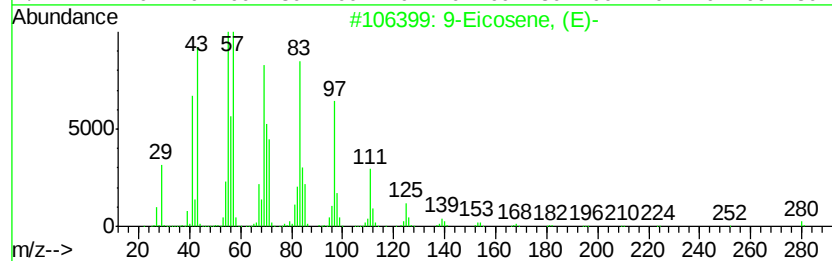
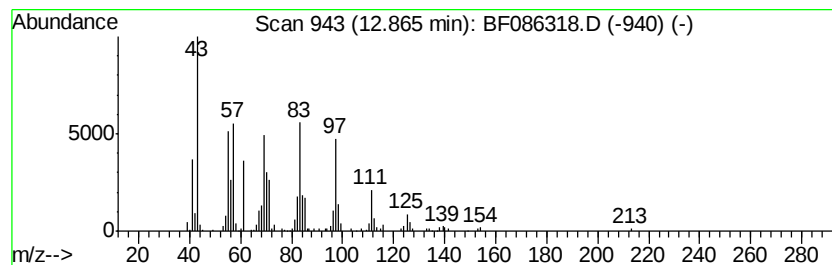
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 9-Eicosene, (E)- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.87	4.70 ng	333987	Chrysene-d12	14.01

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			9-Eicosene, (E)-	280	C20H40	074685-29-3	87
2			5-Octadecene, (E)-	252	C18H36	007206-21-5	87
3			4-Heptafluorobutylrvloxyhexadecane	438	C20H33F7O2	1000282-97-2	86
4			Chloroacetic acid, pentadecyl ester	304	C17H33ClO2	070301-47-2	81
5			1-Heptadecene	238	C17H34	006765-39-5	81



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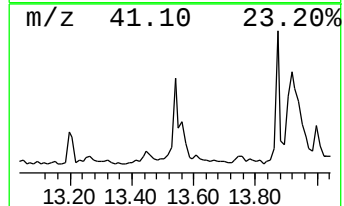
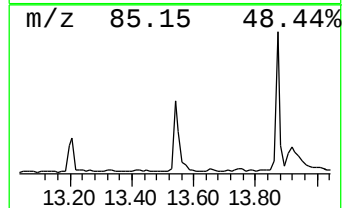
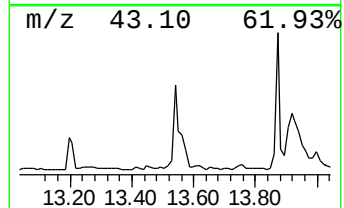
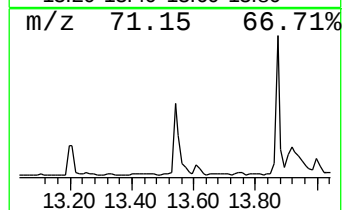
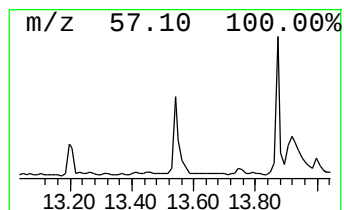
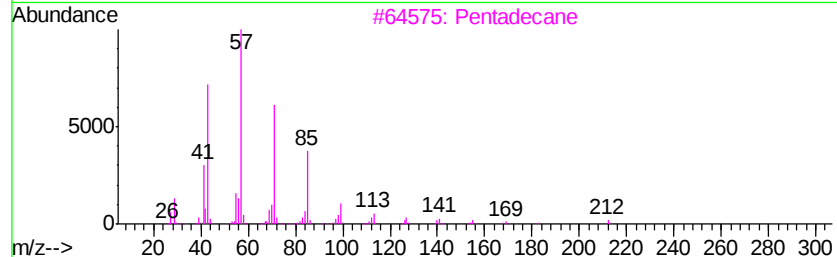
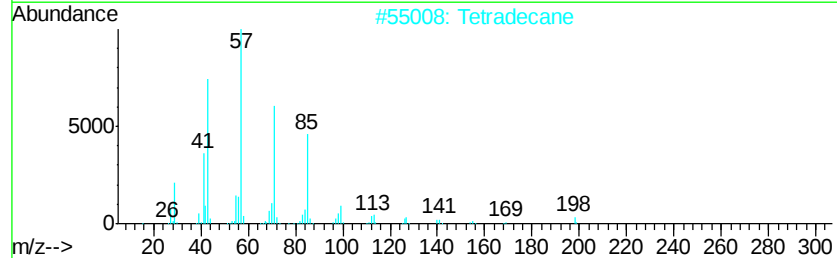
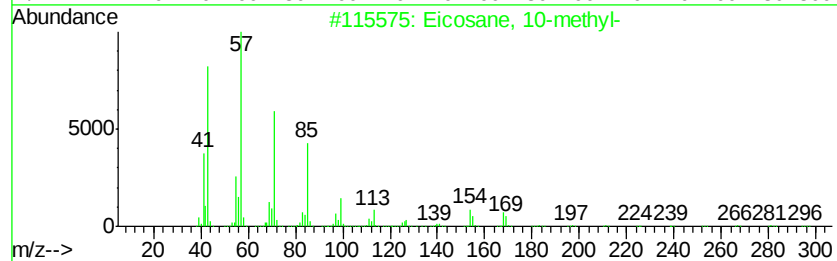
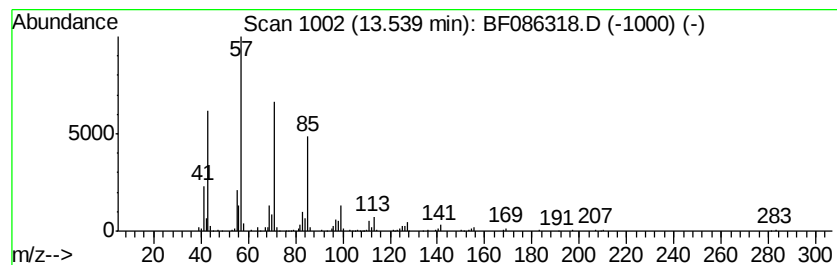
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Peak Number 12 Eicosane, 10-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.54	5.25 ng	373123	Chrysene-d12	14.01

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane, 10-methyl-	296	C21H44	054833-23-7	90
2		Tetradecane	198	C14H30	000629-59-4	86
3		Pentadecane	212	C15H32	000629-62-9	86
4		Octacosane	394	C28H58	000630-02-4	86
5		Heptacosane	380	C27H56	000593-49-7	74



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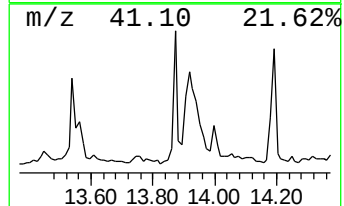
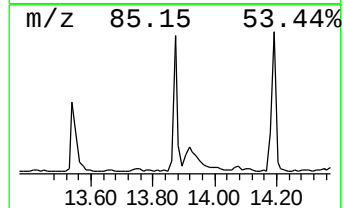
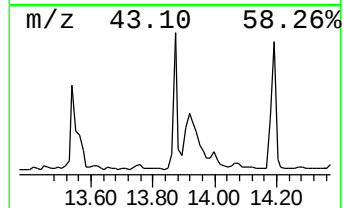
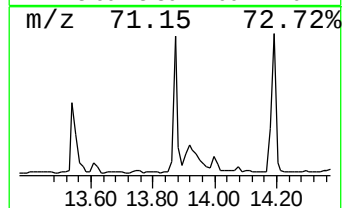
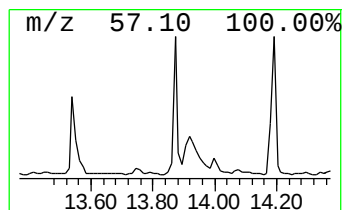
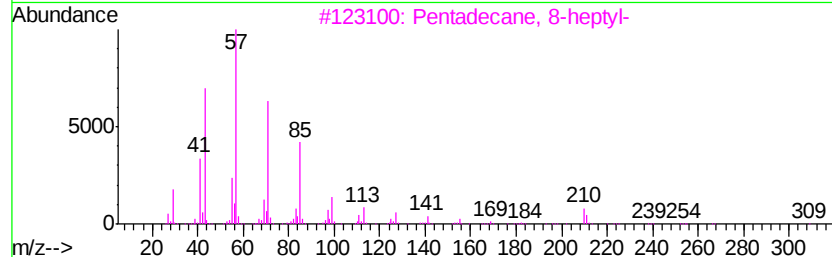
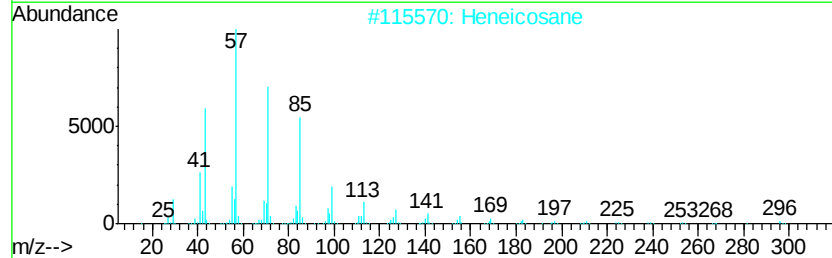
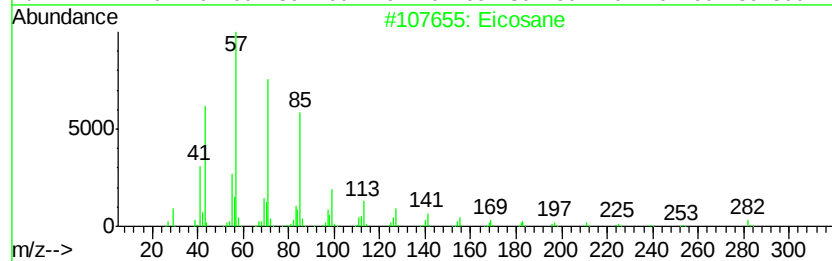
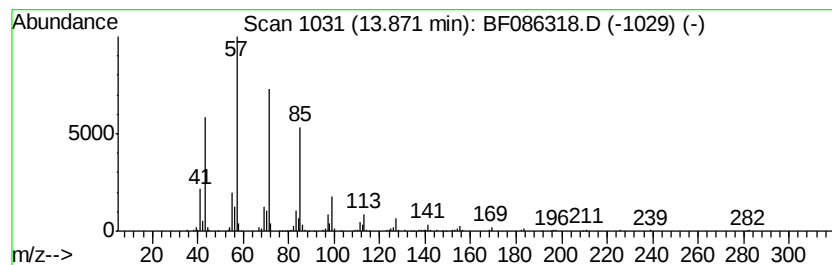
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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Peak Number 13 Eicosane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD		R.T.
13.87	4.98 ng	354051	Chrysene-d12		14.01
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane	282	C20H42	000112-95-8	97
2	Heneicosane	296	C21H44	000629-94-7	91
3	Pentadecane, 8-heptvl-	310	C22H46	071005-15-7	91
4	Nonadecane	268	C19H40	000629-92-5	90
5	Hexadecane	226	C16H34	000544-76-3	90



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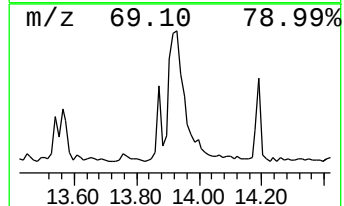
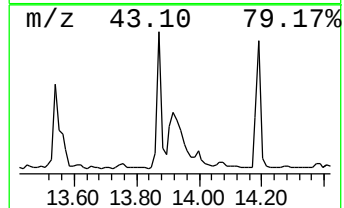
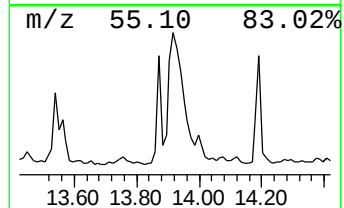
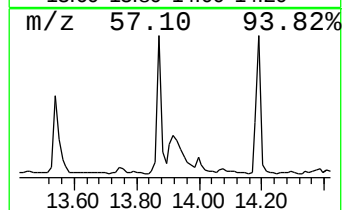
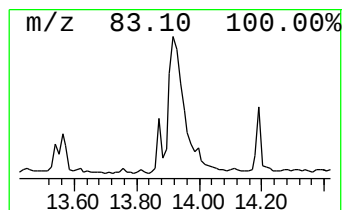
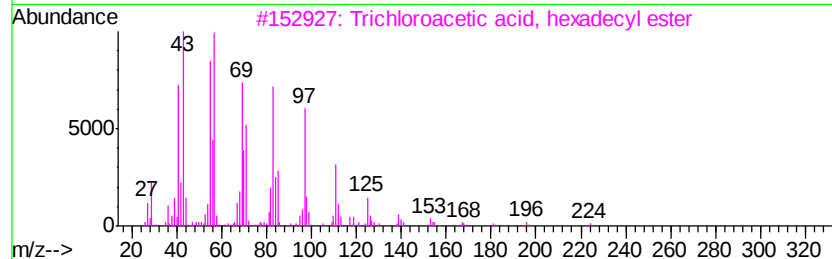
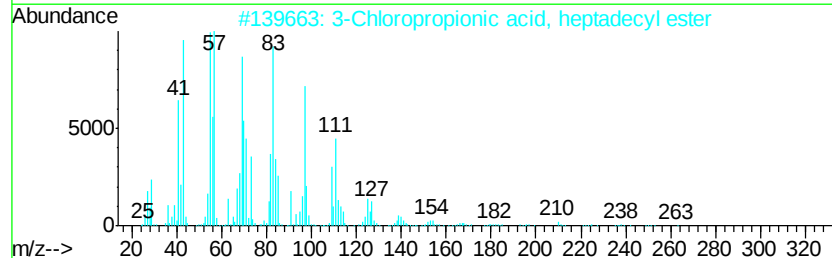
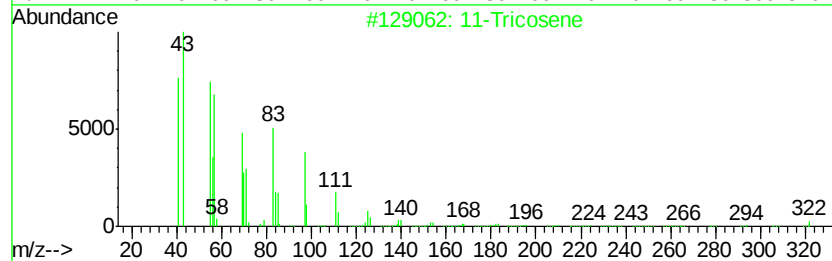
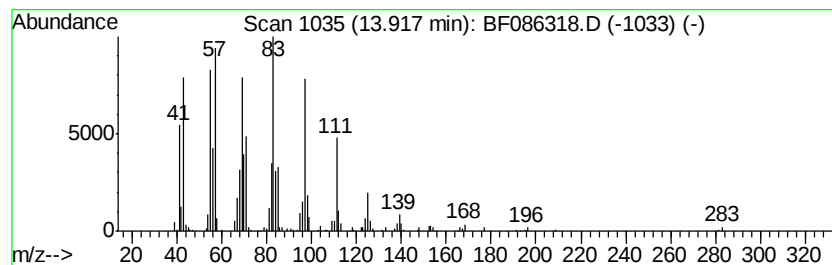
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 11-Tricosene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.92	9.99 ng	709627	Chrysene-d12	14.01

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	11-Tricosene	322	C23H46	052078-56-5	91
2		3-Chloropropionic acid, heptadec...	346	C20H39ClO2	1000283-05-1	91
3		Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	90
4		9-Eicosene, (E)-	280	C20H40	074685-29-3	90
5		Cyclohexadecane	224	C16H32	000295-65-8	87



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF042016\
Data File : BF086318.D
Acq On : 20 Apr 2016 19:43
Operator : UM/SJ
Sample : H2625-05
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MW-2

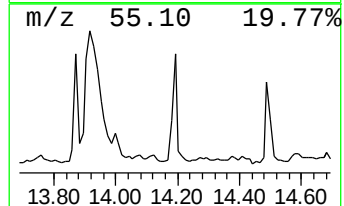
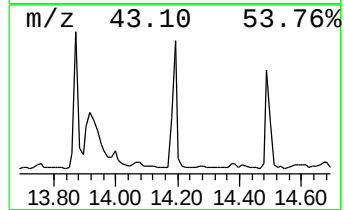
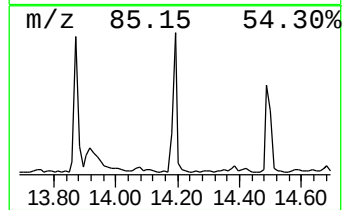
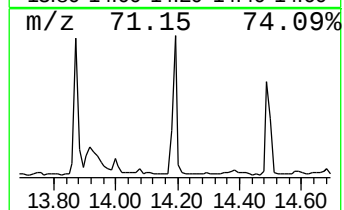
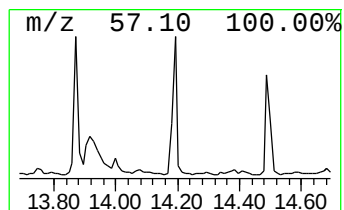
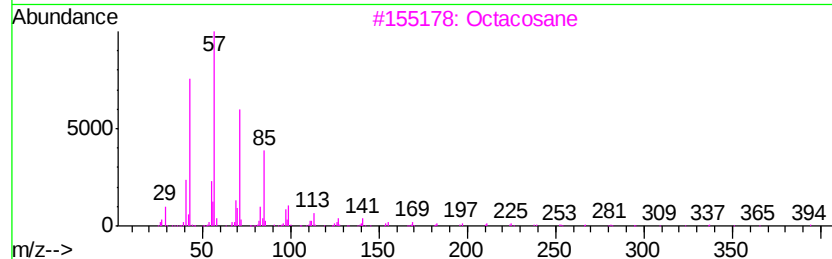
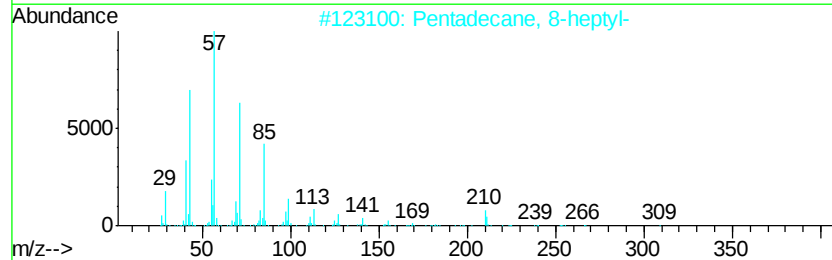
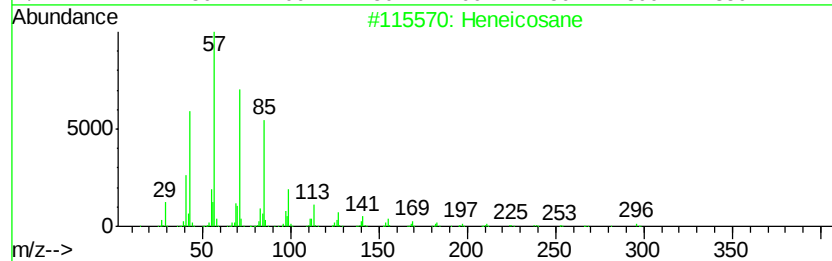
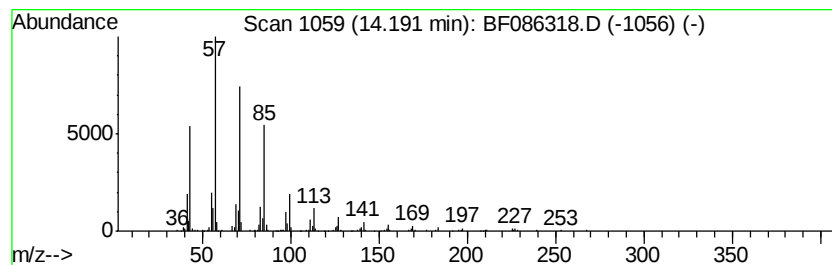
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Heneicosane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.19	5.40 ng	383394	Chrysene-d12	14.01

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	91
2		Pentadecane, 8-heptyl-	310	C22H46	071005-15-7	90
3		Octacosane	394	C28H58	000630-02-4	87
4		Tridecane, 6-propyl-	226	C16H34	055045-10-8	87
5		Tetracosane	338	C24H50	000646-31-1	87



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Data File : BF086318.D
Acq On : 20 Apr 2016 19:43
Operator : UM/SJ
Sample : H2625-05
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MW-2

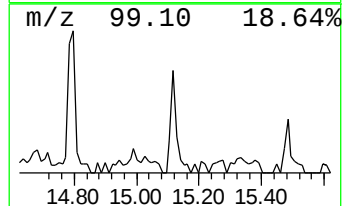
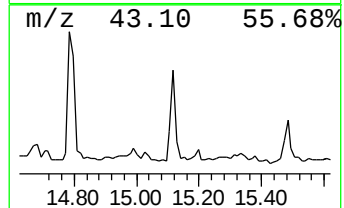
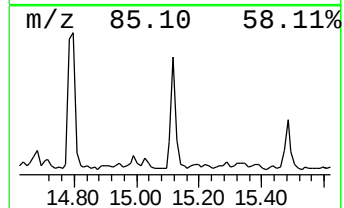
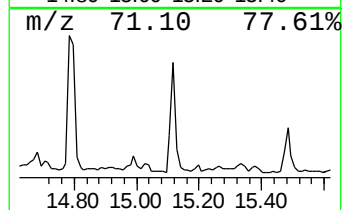
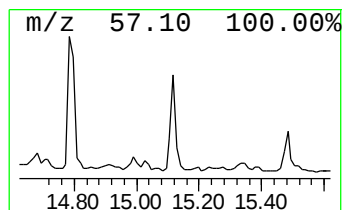
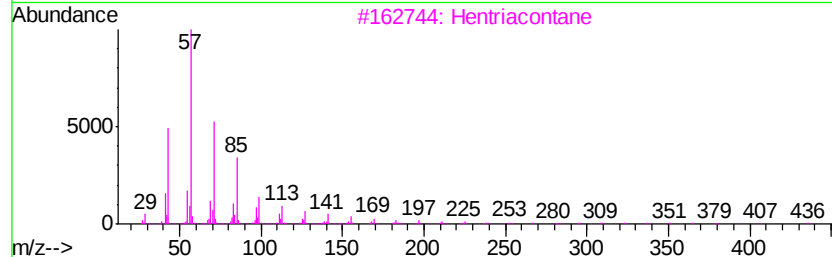
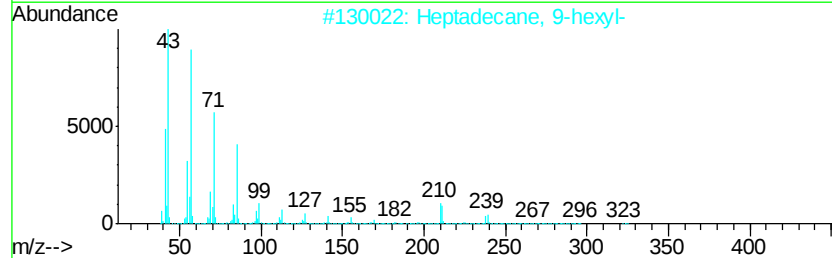
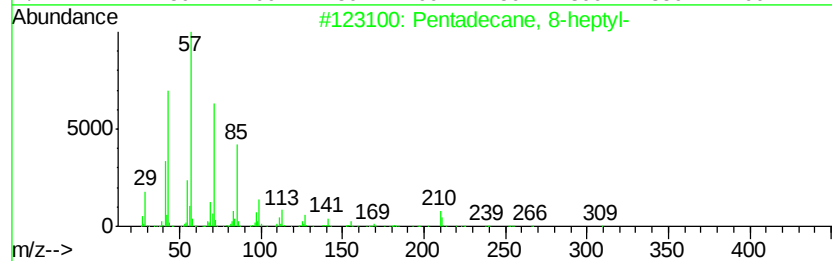
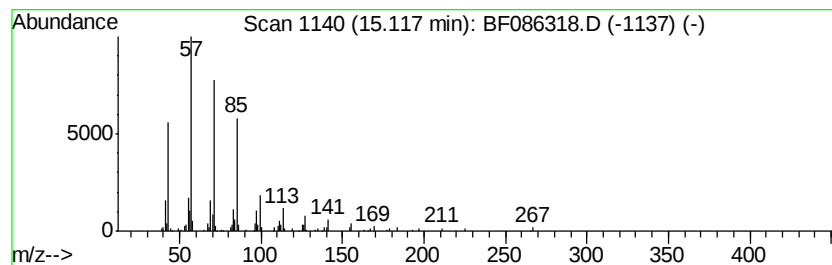
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 Pentadecane, 8-heptyl- Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.12	2.13 ng	126520	Perylene-d12	15.44

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentadecane, 8-heptyl-	310	C22H46	071005-15-7	91
2		Heptadecane, 9-hexyl-	324	C23H48	055124-79-3	90
3		Hentriacontane	437	C31H64	000630-04-6	83
4		Heptacosane	380	C27H56	000593-49-7	83
5		Eicosane, 2-methyl-	296	C21H44	001560-84-5	80



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Sample : H2625-05
Misc :
ALS Vial : 22 Sample Multiplier: 1

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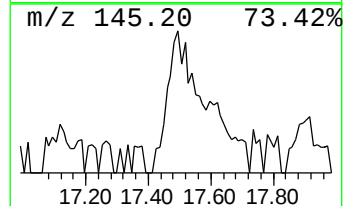
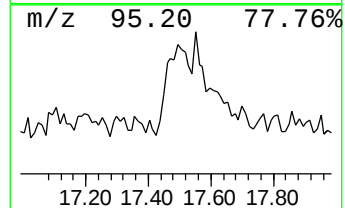
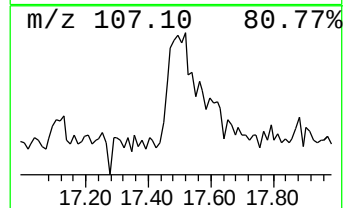
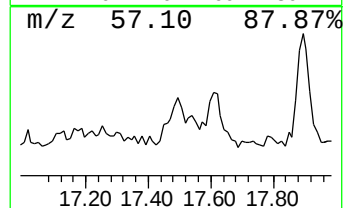
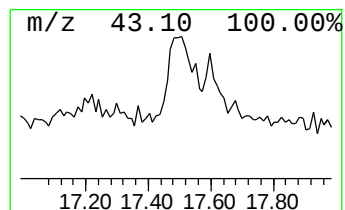
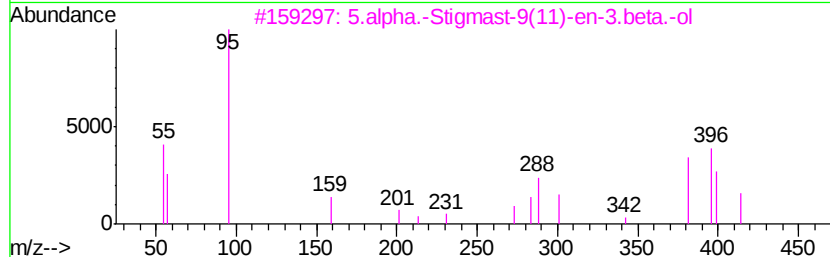
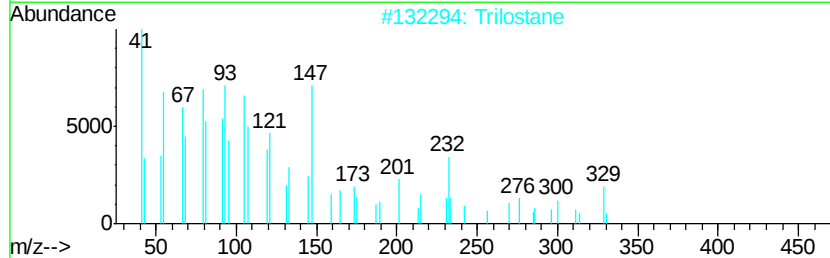
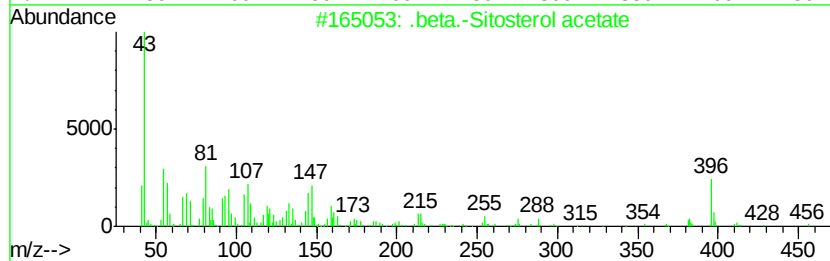
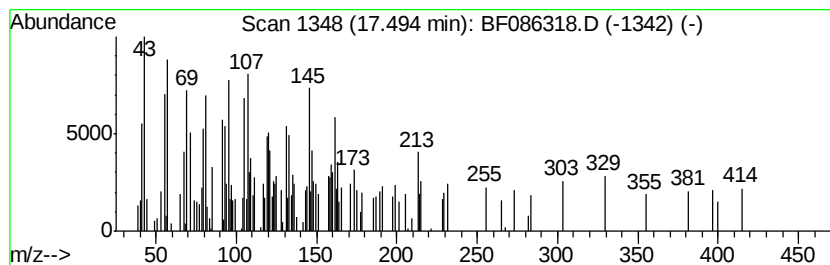
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 19 .beta.-Sitosterol acetate Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.49	4.35 ng	258700	Perylene-d12	15.44

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			.beta.-Sitosterol acetate	456	C31H52O2	000915-05-9	50
2			Trilostane	329	C20H27NO3	013647-35-3	20
3			5.alpha.-Stigmast-9(11)-en-3.beta...	414	C29H50O	077794-81-1	12
4			5.beta.,6.beta.-Epoxy-7-bromocho...	478	C27H43BrO2	104825-82-3	10
5			N-Nitrososolasodine	442	C27H42N2O3	004860-12-2	10



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 Operator : UM/SJ
 Sample : H2625-05
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
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 ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF042016.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
Propane, 1,1-dime...	2.18	13.4	ng	914282	1	6.86	1367930	20.0
2-Butanol, 2,3-di...	3.06	2.3	ng	154603	1	6.86	1367930	20.0
2-Pentanone, 4-hy...	5.05	19.5	ng	1331460	1	6.86	1367930	20.0
2-Pentanone, 4-me...	5.87	3.6	ng	245782	1	6.86	1367930	20.0
unknown6.62	6.62	66.9	ng	4573610	1	6.86	1367930	20.0
unknown7.50	7.50	2.3	ng	233390	2	8.14	2045520	20.0
Tetradecane	10.34	3.8	ng	388348	3	9.89	2031270	20.0
Decane, 2,3,5-tri...	10.81	4.7	ng	432438	4	11.38	1841260	20.0
9-Eicosene, (E)-	12.87	4.7	ng	333987	5	14.01	1420730	20.0
Eicosane, 10-methyl-	13.54	5.3	ng	373123	5	14.01	1420730	20.0
Eicosane	13.87	5.0	ng	354051	5	14.01	1420730	20.0
11-Tricosene	13.92	10.0	ng	709627	5	14.01	1420730	20.0
Heneicosane	14.19	5.4	ng	383394	5	14.01	1420730	20.0
Pentadecane, 8-he...	15.12	2.1	ng	126520	6	15.44	1190670	20.0
.beta.-Sitosterol...	17.49	4.3	ng	258700	6	15.44	1190670	20.0