

Data Path : Z:\HPCHEM1\BNA F\DATA\BF042716\
 Data File : BF086460.D
 Acq On : 28 Apr 2016 8:33
 Operator : UM/SJ
 Sample : H2736-04
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RICHMOND-AMBOY-SB01(0.5-10.0)

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-BF042716.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	4.990	252	254	262	rVB	63553	112503	2.82%	0.363%
2	5.402	287	290	293	rBV	3567305	3758843	94.06%	12.117%
3	6.442	378	381	384	rBV	3152382	3703431	92.68%	11.938%
4	6.579	390	393	395	rBV	2893888	3996096	100.00%	12.881%
5	6.807	410	413	415	rBV	704703	731362	18.30%	2.358%
6	6.956	424	426	429	rBV	2077893	2885709	72.21%	9.302%
7	7.367	459	462	465	rBV	1937347	2193451	54.89%	7.071%
8	7.470	469	471	479	rVV	54939	153693	3.85%	0.495%
9	8.087	523	525	528	rBV	931668	871415	21.81%	2.809%
10	9.173	617	620	622	rBV	2411091	3216128	80.48%	10.367%
11	9.288	629	630	642	rVB2	18399	40637	1.02%	0.131%
12	9.562	652	654	656	rBV	574650	480968	12.04%	1.550%
13	9.779	671	673	676	rBV	62568	57420	1.44%	0.185%
14	9.848	676	679	681	rBV	490672	705956	17.67%	2.276%
15	10.282	715	717	718	rVB	111548	84873	2.12%	0.274%
16	10.499	733	736	739	rBV	38170	69726	1.74%	0.225%
17	10.625	745	747	750	rBV2	1416474	1708539	42.76%	5.508%
18	10.751	756	758	765	rVB	125579	158430	3.96%	0.511%
19	11.196	796	797	799	rBV	60258	61368	1.54%	0.198%
20	11.322	806	808	812	rBV	615191	623763	15.61%	2.011%
21	12.911	944	947	949	rBV	2636785	2608302	65.27%	8.408%
22	13.402	987	990	992	rBV	19127	41559	1.04%	0.134%
23	13.814	1023	1026	1034	rBV	194141	350531	8.77%	1.130%
24	13.951	1036	1038	1046	rVV	619521	806792	20.19%	2.601%
25	14.065	1046	1048	1050	rVB	94567	106147	2.66%	0.342%
26	14.454	1077	1082	1087	rBV4	42565	157260	3.94%	0.507%
27	14.797	1110	1112	1115	rVB	56764	59597	1.49%	0.192%
28	15.048	1132	1134	1136	rBV	58119	60274	1.51%	0.194%
29	15.357	1159	1161	1164	rBV	311475	515882	12.91%	1.663%
30	15.814	1198	1201	1204	rBV	43738	75507	1.89%	0.243%
31	17.471	1343	1346	1353	rVB3	23092	66883	1.67%	0.216%
32	17.780	1371	1373	1379	rVB6	21915	57755	1.45%	0.186%
33	18.031	1391	1395	1403	rVB7	20359	79495	1.99%	0.256%
34	18.489	1431	1435	1440	rBV7	15237	51068	1.28%	0.165%

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35	19.174	1490	1495	1502	rVB3	113773	370617	9.27%	1.195%
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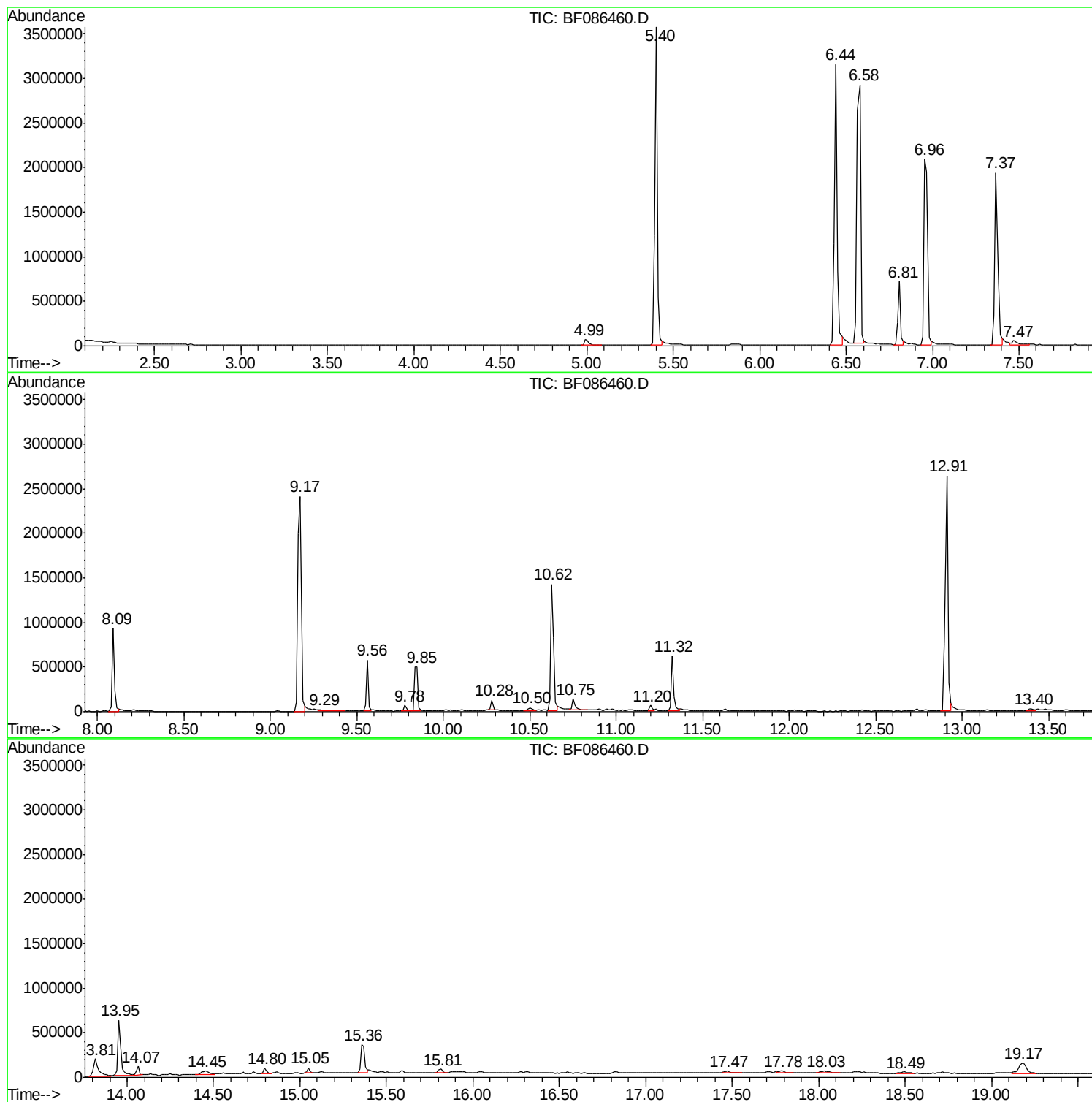
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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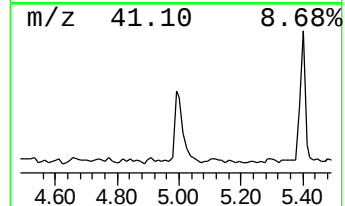
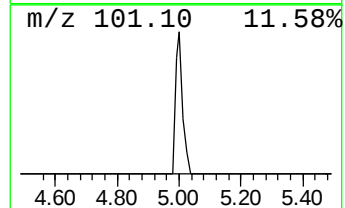
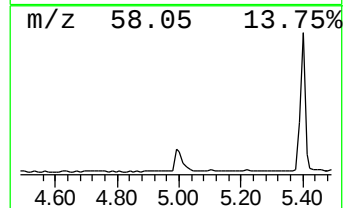
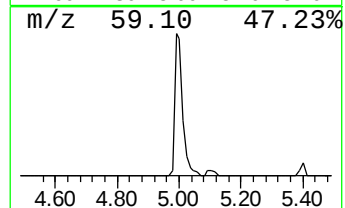
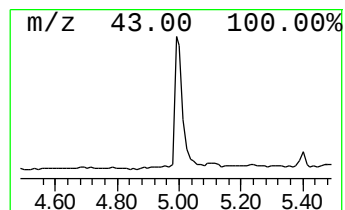
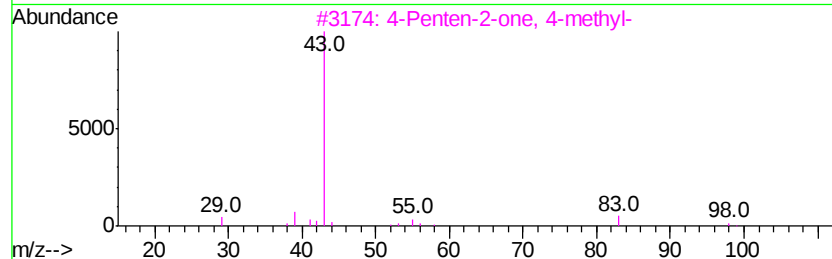
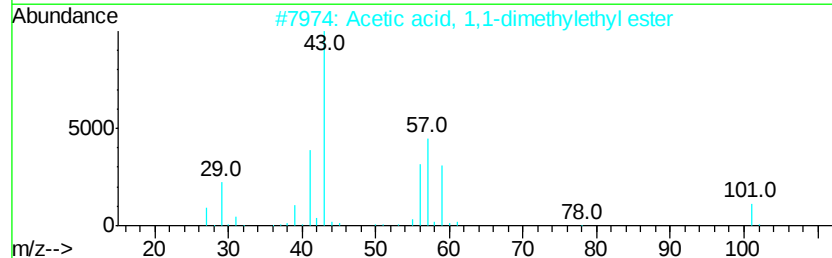
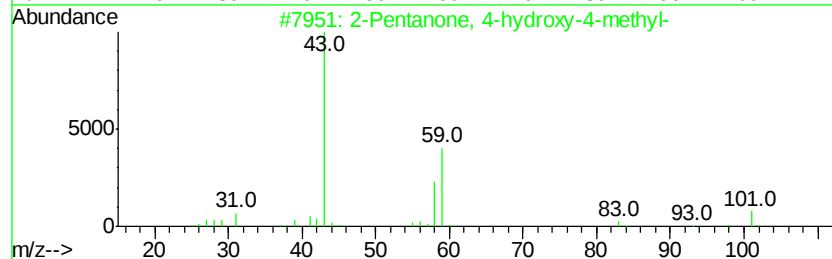
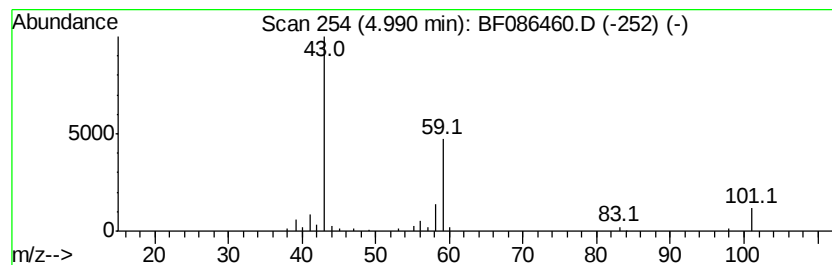
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.99	3.08 ng	112503	1,4-Dichlorobenzene-d4	6.81

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	28
3			4-Penten-2-one, 4-methyl-	98	C6H10O	003744-02-3	14
4			Diacetamide	101	C4H7NO2	000625-77-4	9
5			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9



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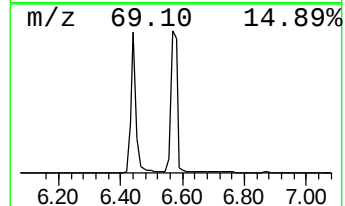
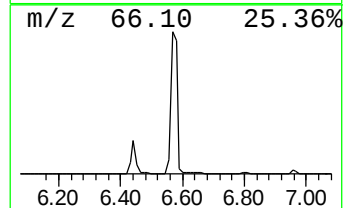
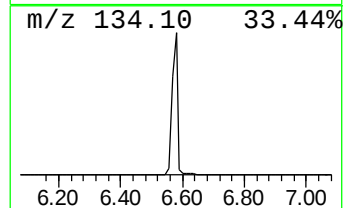
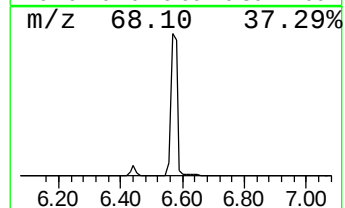
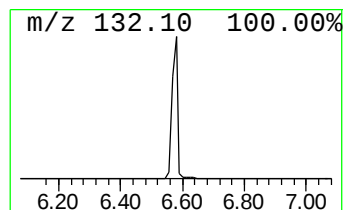
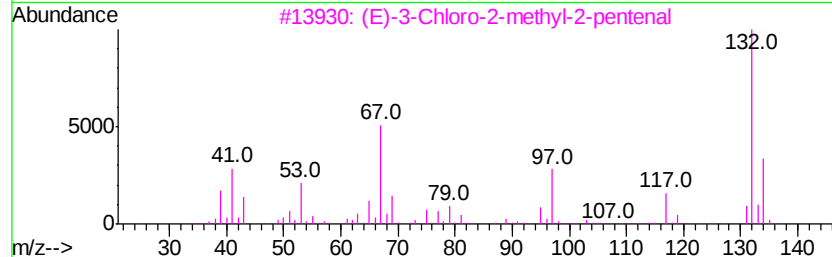
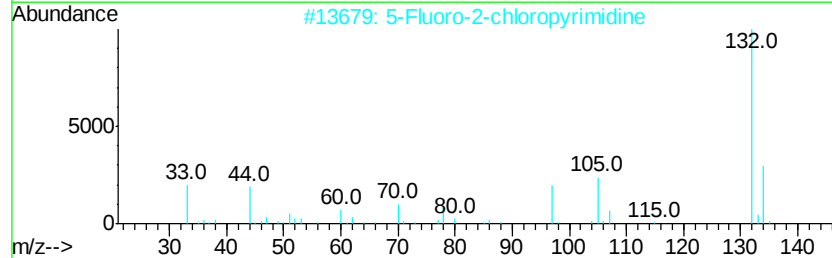
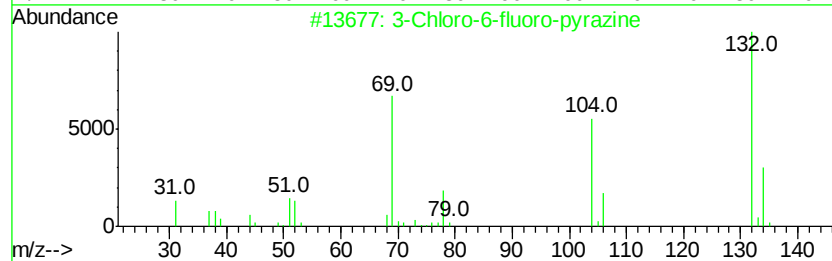
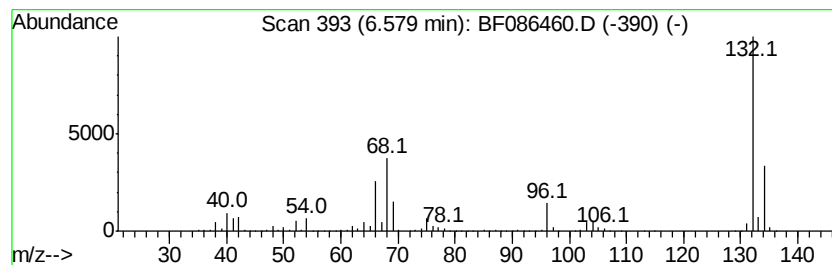
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown6.58 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.58	109.28 ng	3996100	1,4-Dichlorobenzene-d4	6.81

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	35
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	25
3		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	16
5		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9



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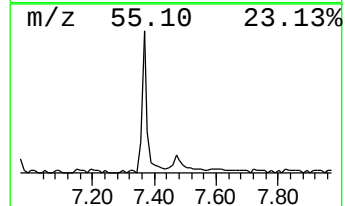
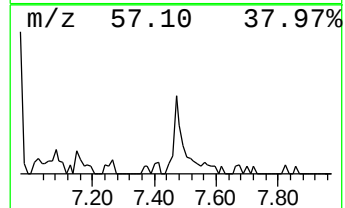
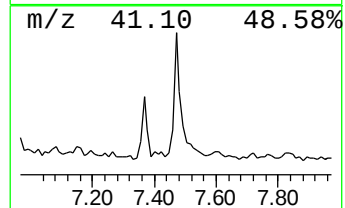
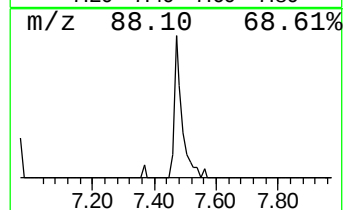
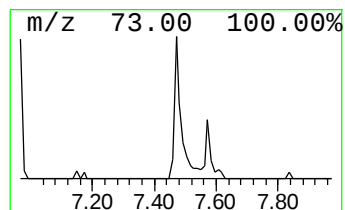
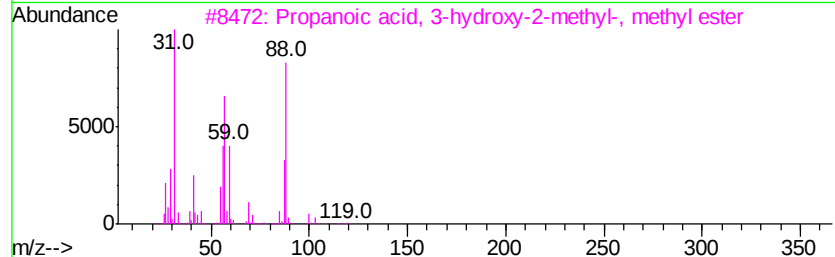
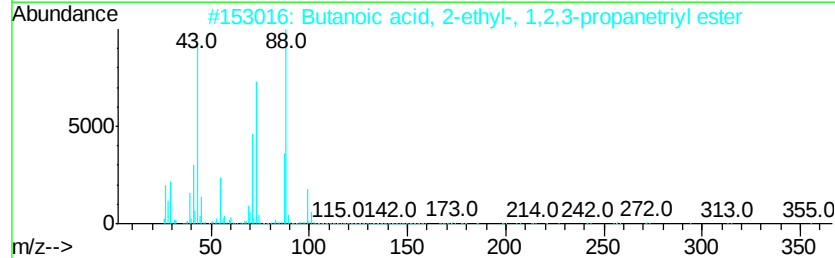
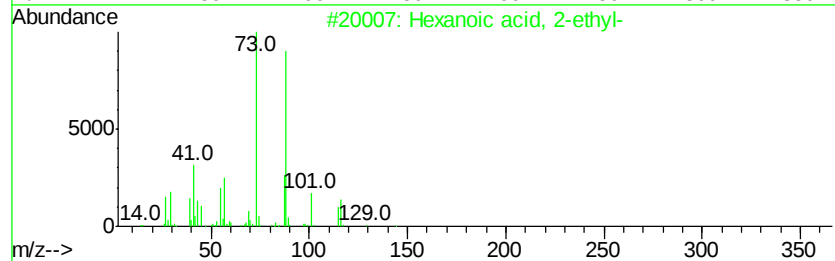
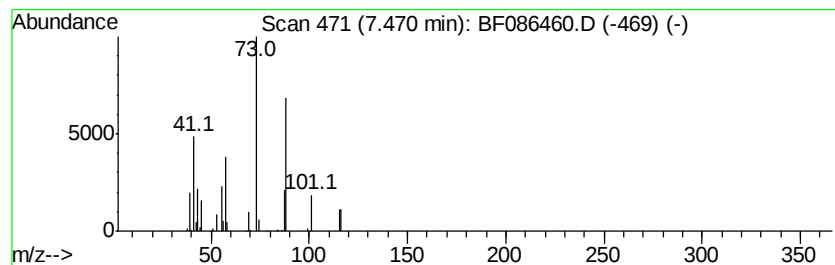
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Peak Number 4 Hexanoic acid, 2-ethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.47	3.53 ng	153693	Naphthalene-d8	8.09

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	83
2			Butanoic acid, 2-ethyl-, 1,2,3-p...	386	C21H38O6	056554-54-2	40
3			Propanoic acid, 3-hydroxy-2-meth...	118	C5H10O3	042998-03-8	23
4			3-Hexanol, 3,5-dimethyl-	130	C8H18O	004209-91-0	16
5			Hydrazine, 1-butyl-1-ethyl-	116	C6H16N2	052728-56-0	12



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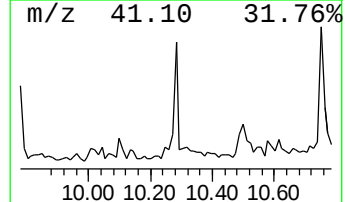
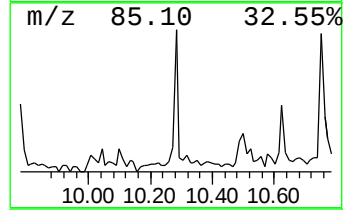
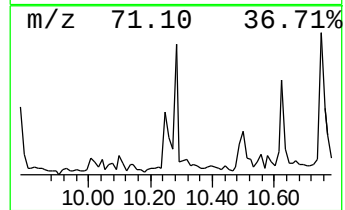
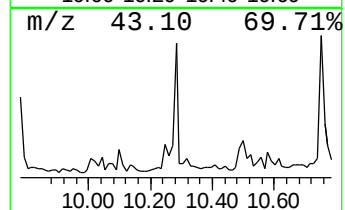
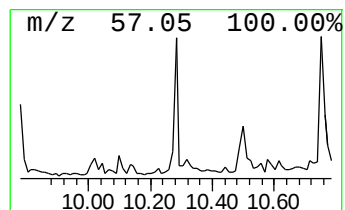
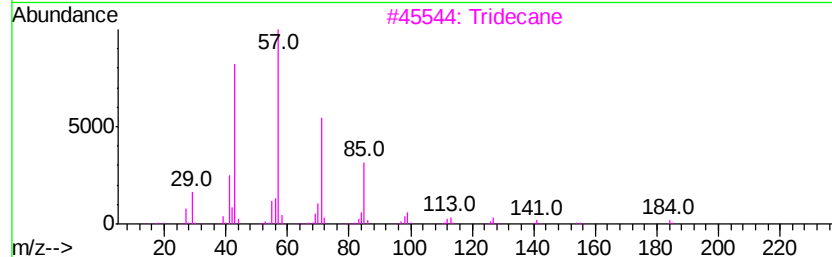
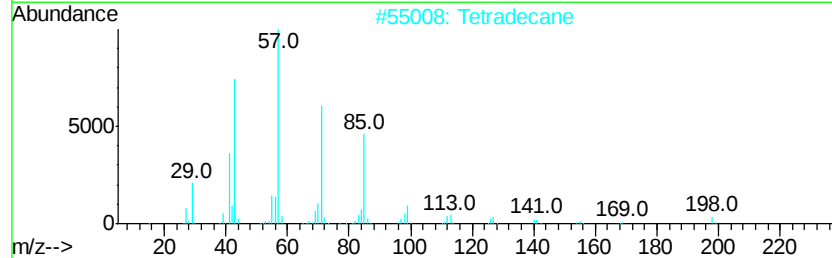
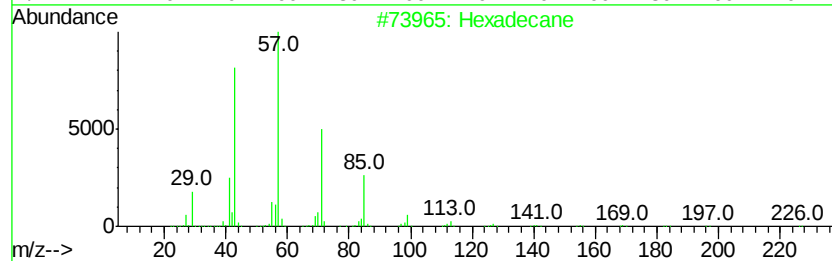
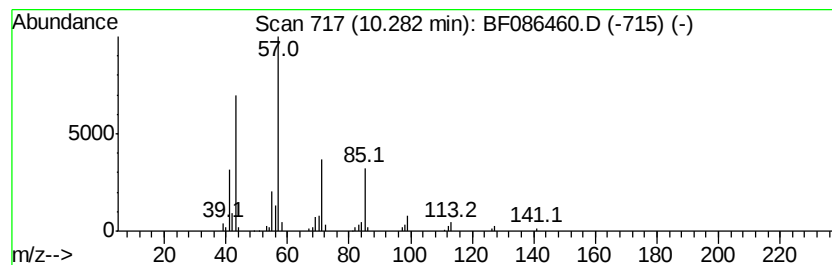
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Peak Number 5 Hexadecane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD		R.T.
10.28	2.40 ng	84873	Acenaphthene-d10		9.85
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane	226	C16H34	000544-76-3	90
2	Tetradecane	198	C14H30	000629-59-4	86
3	Tridecane	184	C13H28	000629-50-5	86
4	Octane, 2,4,6-trimethyl-	156	C11H24	062016-37-9	64
5	Silane, trichlorooctadecyl-	386	C18H37Cl3Si	000112-04-9	64



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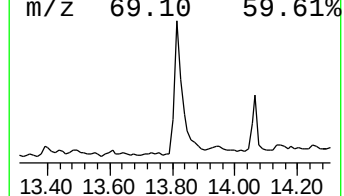
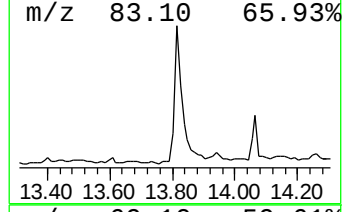
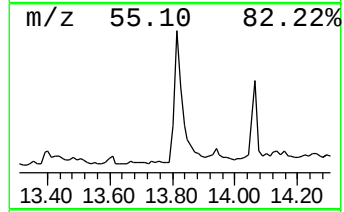
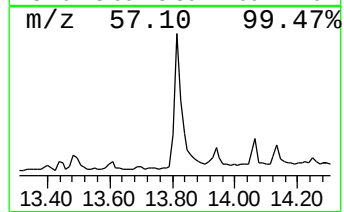
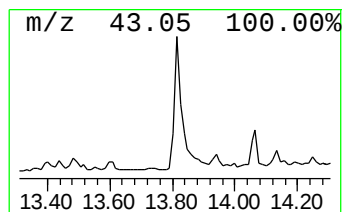
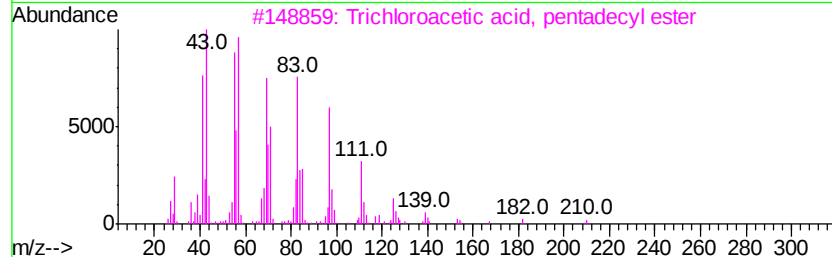
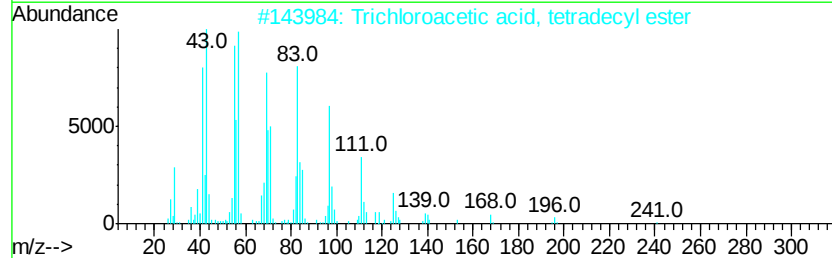
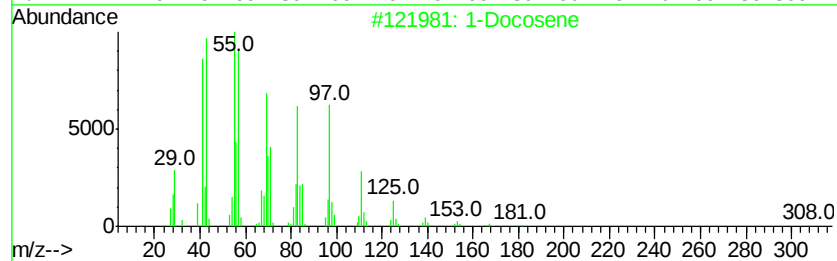
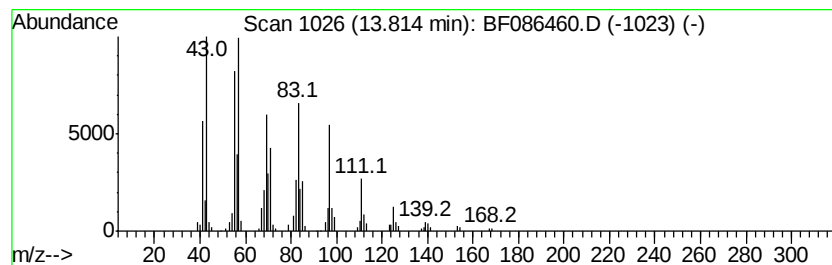
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Peak Number 7 1-Docosene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.81	8.69 ng	350531	Chrysene-d12	13.95

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Docosene	308	C22H44	001599-67-3	95
2			Trichloroacetic acid, tetradecyl...	358	C16H29Cl3O2	074339-52-9	91
3			Trichloroacetic acid, pentadecyl...	372	C17H31Cl3O2	074339-53-0	91
4			Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	91
5			1-Heptadecanol	256	C17H36O	001454-85-9	87



Data Path : Z:\HPCHEM1\BNA F\DATA\BF042716\
Data File : BF086460.D
Acq On : 28 Apr 2016 8:33
Operator : UM/SJ
Sample : H2736-04
Misc :
ALS Vial : 34 Sample Multiplier: 1

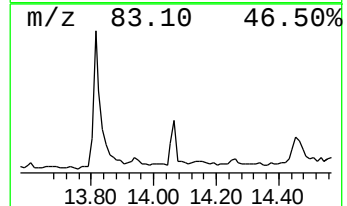
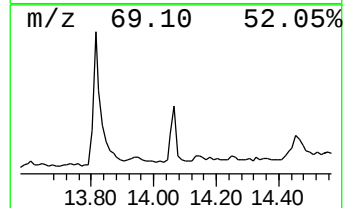
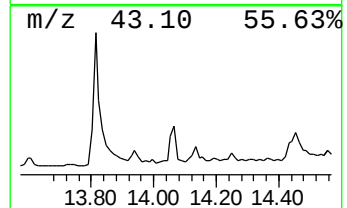
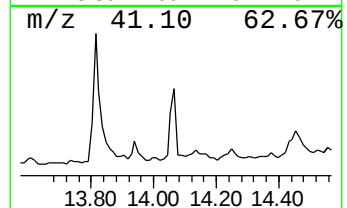
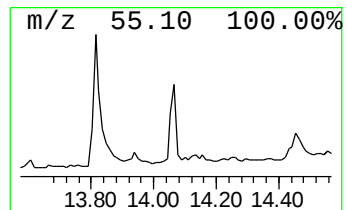
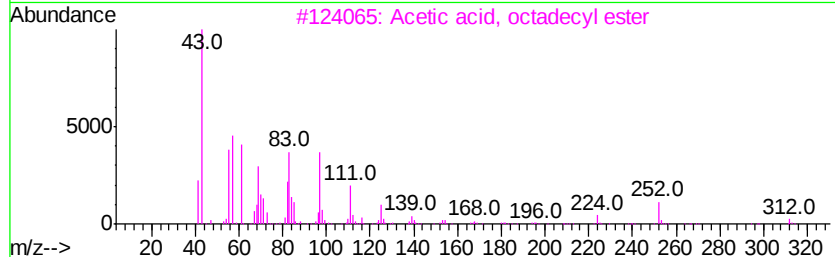
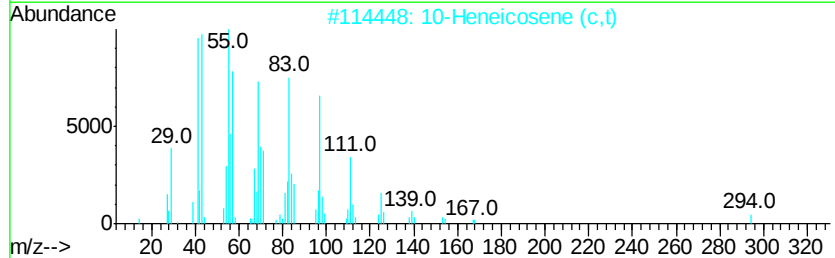
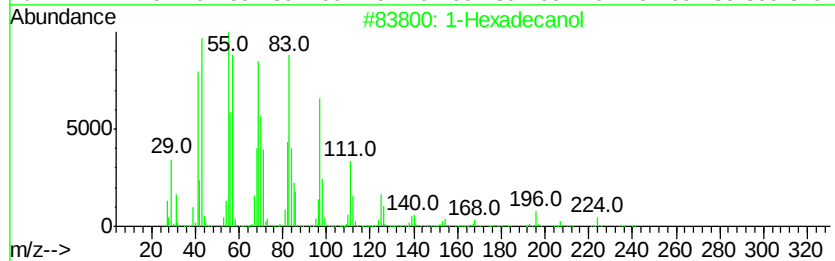
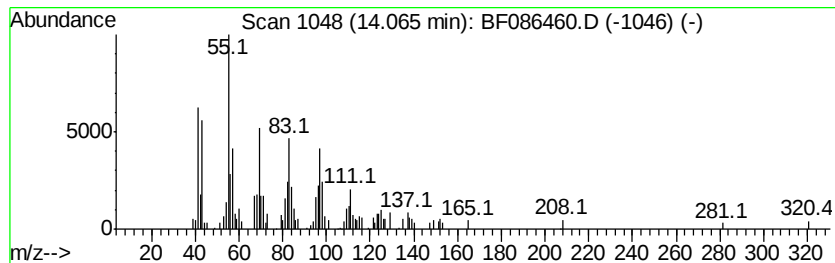
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ClientSampleId :
RICHMOND-AMBOY-SB01(0.5-10.0)

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF042716.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 1-Hexadecanol Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.07	2.63 ng	106147	Chrysene-d12	13.95	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Hexadecanol	242	C16H34O	036653-82-4	64
2	10-Heneicosene (c,t)	294	C21H42	095008-11-0	62
3	Acetic acid, octadecyl ester	312	C20H40O2	000822-23-1	58
4	Chloromethyl 9-chloroundecanoate	268	C12H22Cl2O2	080418-95-7	58
5	1-Octadecanol	270	C18H38O	000112-92-5	58



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF042716\
Data File : BF086460.D
Acq On : 28 Apr 2016 8:33
Operator : UM/SJ
Sample : H2736-04
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
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ClientSampleId :
RICHMOND-AMBOY-SB01(0.5-10.0)

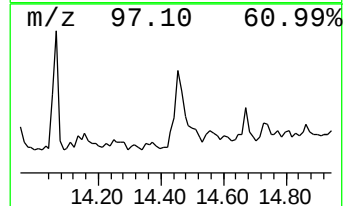
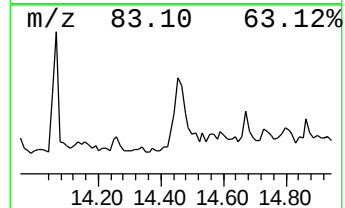
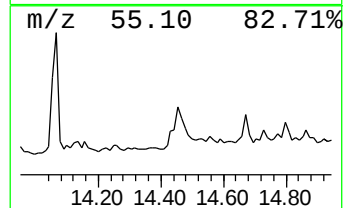
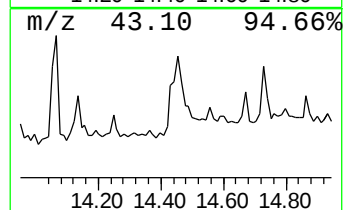
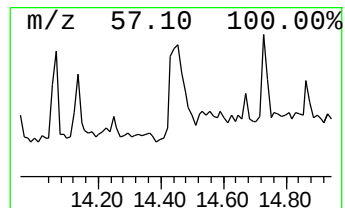
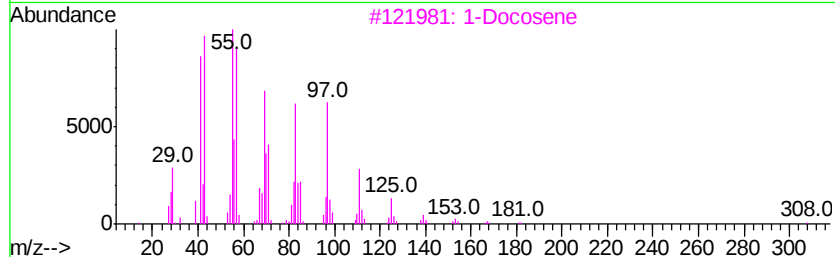
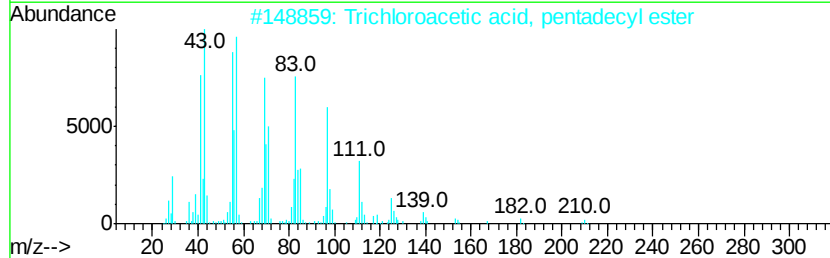
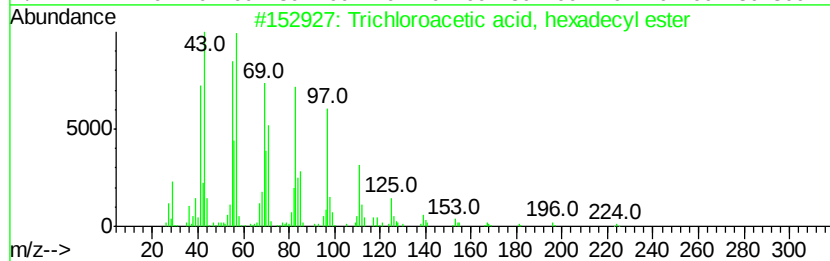
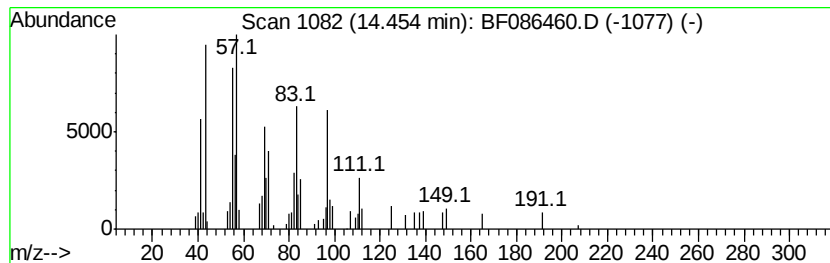
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Trichloroacetic acid, hexad... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.45	3.90 ng	157260	Chrysene-d12	13.95

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	90
2			Trichloroacetic acid, pentadecyl...	372	C17H31Cl3O2	074339-53-0	90
3			1-Docosene	308	C22H44	001599-67-3	87
4			Trifluoroacetic acid, n-heptadec...	324	C17H31F3O2	1000216-79-2	86
5			17-Pentatriacontene	491	C35H70	006971-40-0	80



Data Path : Z:\HPCHEM1\BNA F\DATA\BF042716\
Data File : BF086460.D
Acq On : 28 Apr 2016 8:33
Operator : UM/SJ
Sample : H2736-04
Misc :
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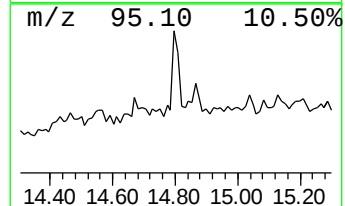
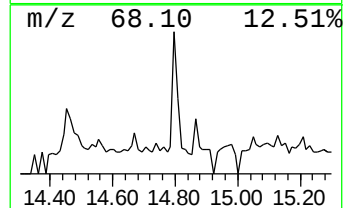
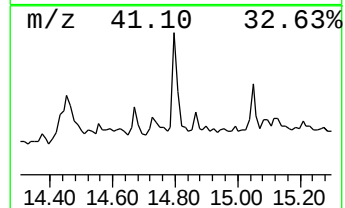
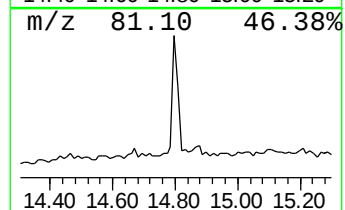
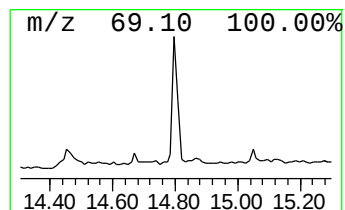
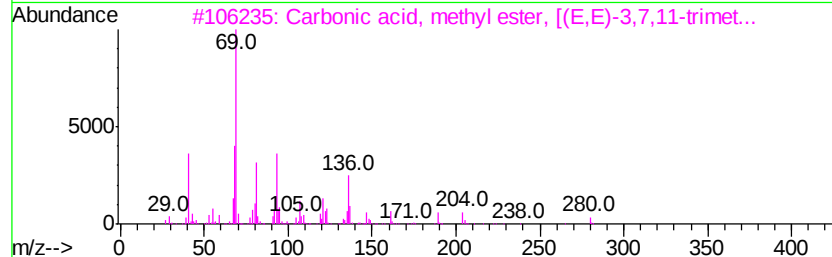
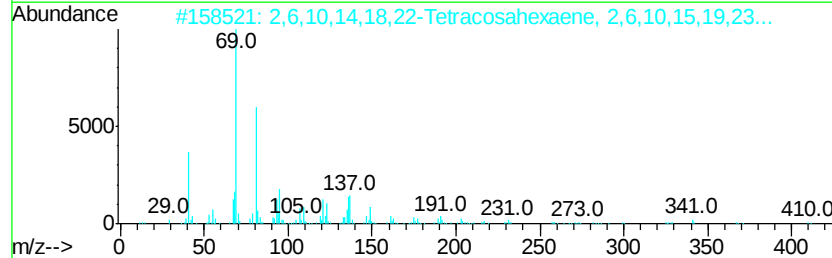
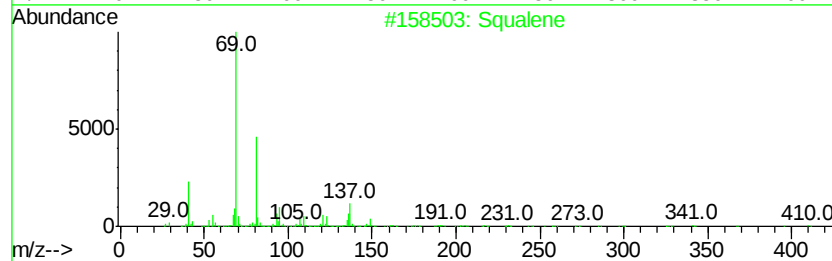
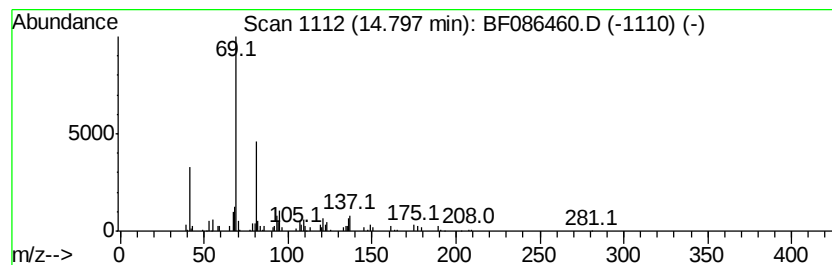
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Squalene Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.80	2.31 ng	59597	Perylene-d12	15.37

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Squalene	410	C30H50	007683-64-9	90
2		2,6,10,14,18,22-Tetracosahexaene...	410	C30H50	000111-02-4	83
3		Carbonic acid, methyl ester, [(E...]	280	C17H28O3	1000164-38-7	72
4		2,6,10-Dodecatrien-1-ol, 3,7,11-...	222	C15H26O	000106-28-5	72
5		7,11-Dimethyldodeca-2,6,10-trien...	208	C14H24O	125529-10-4	72



Data Path : Z:\HPCHEM1\BNA F\DATA\BF042716\
Data File : BF086460.D
Acq On : 28 Apr 2016 8:33
Operator : UM/SJ
Sample : H2736-04
Misc :
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Instrument :
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ClientSampleId :
RICHMOND-AMBOY-SB01(0.5-10.0)

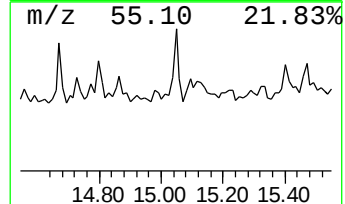
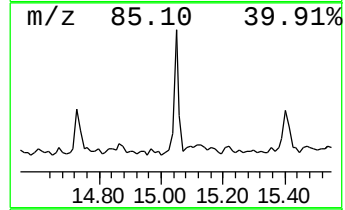
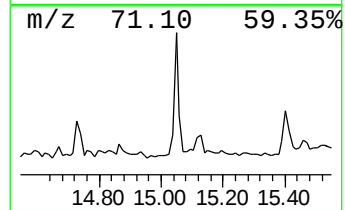
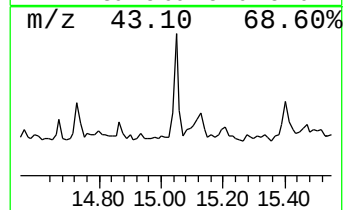
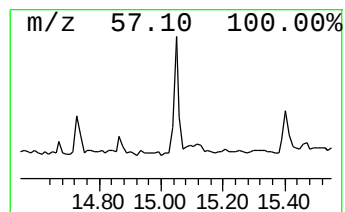
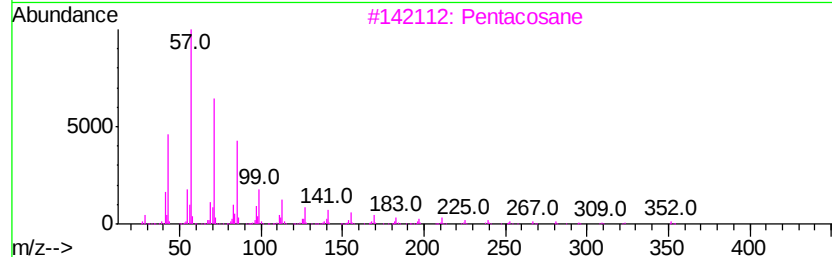
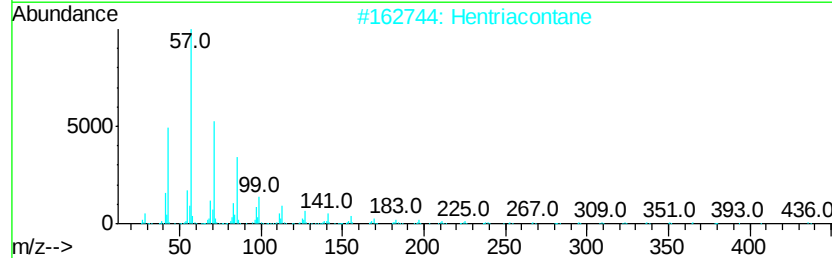
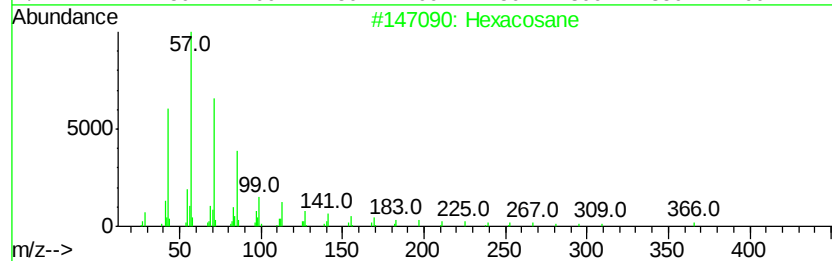
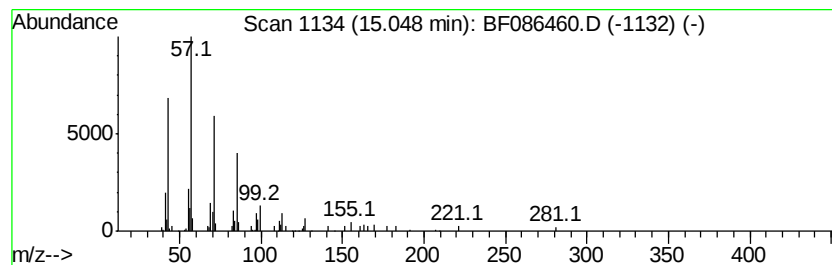
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Hexacosane Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.05	2.34 ng	60274	Perylene-d12	15.37

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexacosane	366	C26H54	000630-01-3	91
2			Hentriacontane	437	C31H64	000630-04-6	91
3			Pentacosane	352	C25H52	000629-99-2	91
4			Octacosane	394	C28H58	000630-02-4	90
5			Triacontane	422	C30H62	000638-68-6	90



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF042716\
Data File : BF086460.D
Acq On : 28 Apr 2016 8:33
Operator : UM/SJ
Sample : H2736-04
Misc :
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Instrument :
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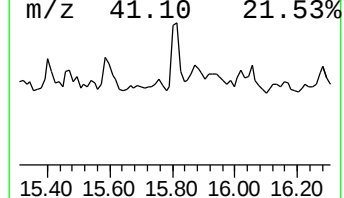
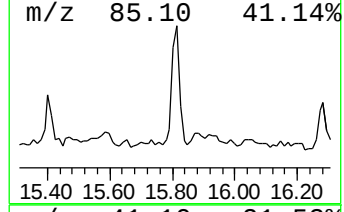
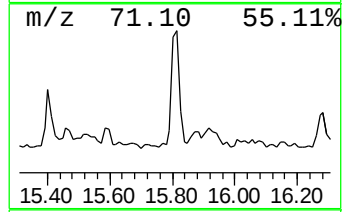
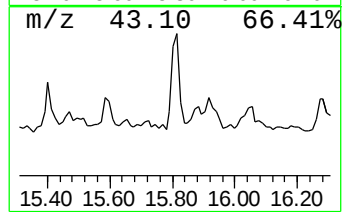
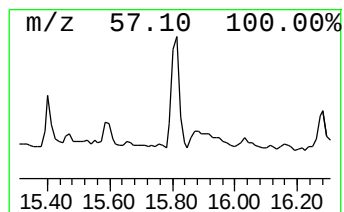
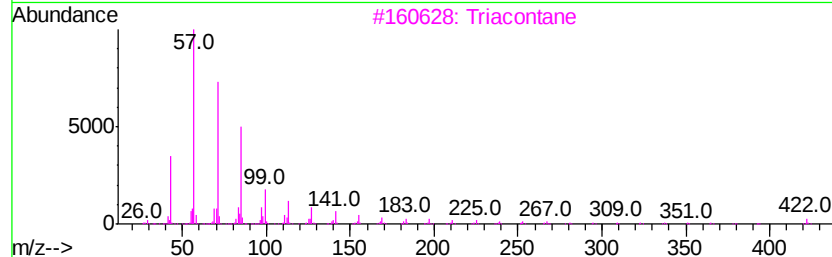
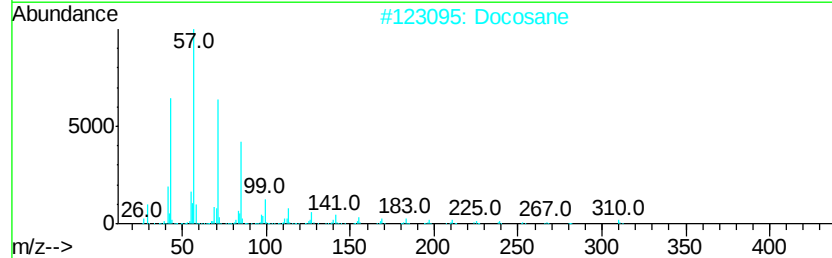
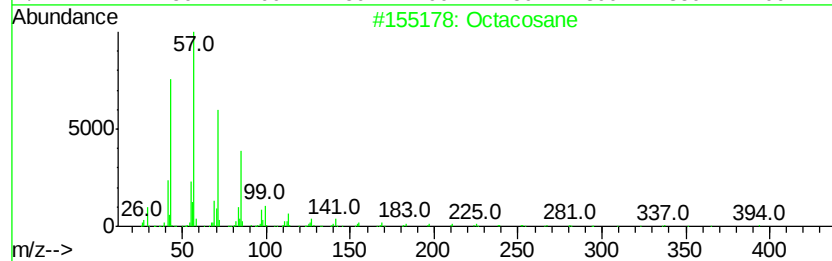
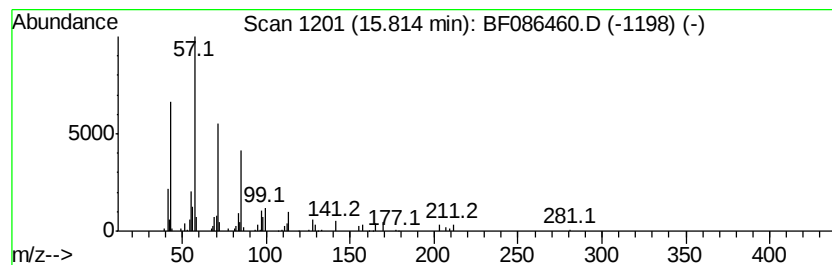
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Octacosane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.81	2.93 ng	75507	Perylene-d12	15.37

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octacosane	394	C28H58	000630-02-4	87
2			Docosane	310	C22H46	000629-97-0	86
3			Triacontane	422	C30H62	000638-68-6	80
4			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	80
5			Heptacosane	380	C27H56	000593-49-7	80



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF042716\
Data File : BF086460.D
Acq On : 28 Apr 2016 8:33
Operator : UM/SJ
Sample : H2736-04
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Instrument :
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ClientSampleId :
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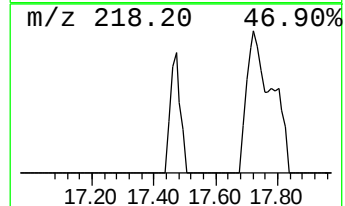
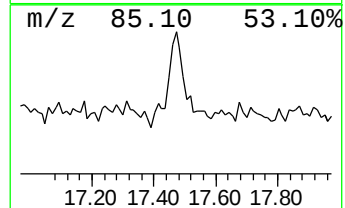
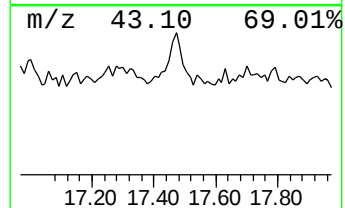
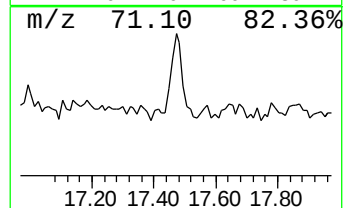
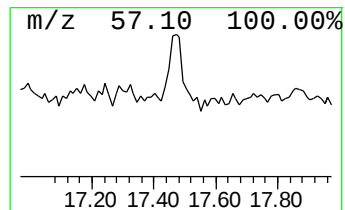
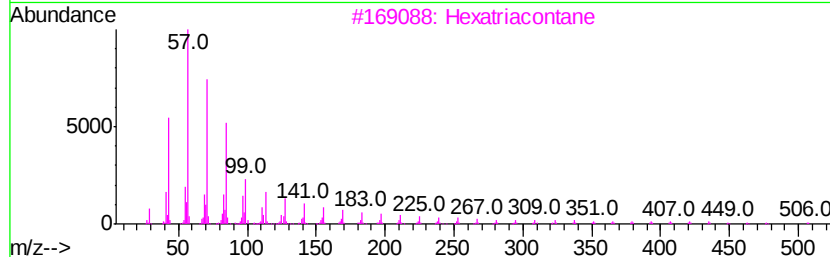
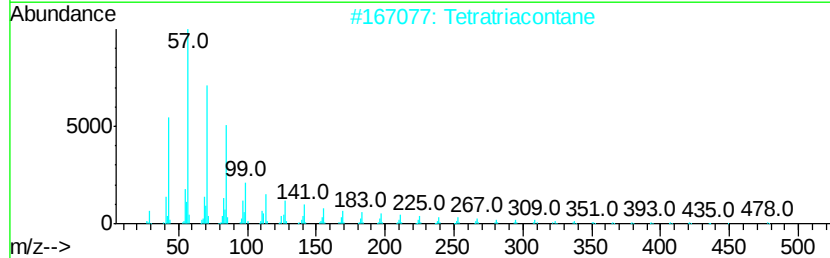
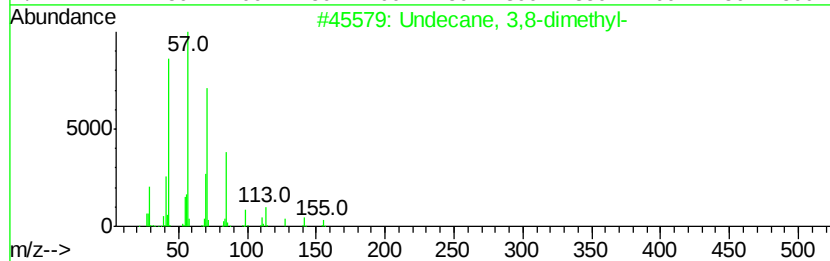
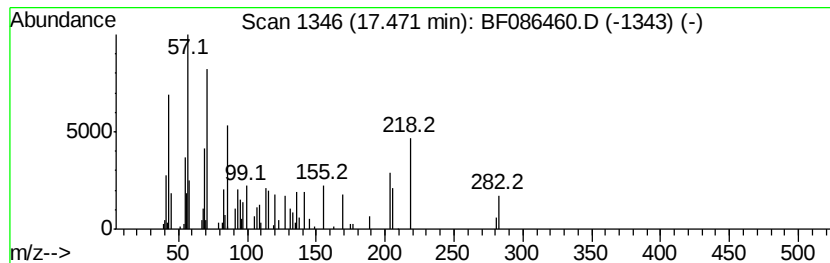
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 unknown17.47 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.47	2.59 ng	66883	Perylene-d12	15.37

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Undecane, 3,8-dimethyl-	184	C13H28	017301-30-3	43
2			Tetratriacontane	479	C34H70	014167-59-0	38
3			Hexatriacontane	507	C36H74	000630-06-8	38
4			Hentriacontane	437	C31H64	000630-04-6	38
5			Hexadecane, 2,6,10,14-tetramethyl-	282	C20H42	000638-36-8	38



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF042716\
Data File : BF086460.D
Acq On : 28 Apr 2016 8:33
Operator : UM/SJ
Sample : H2736-04
Misc :
ALS Vial : 34 Sample Multiplier: 1

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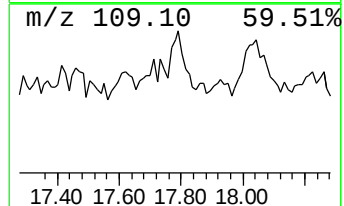
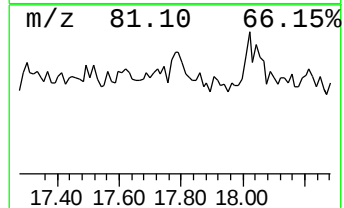
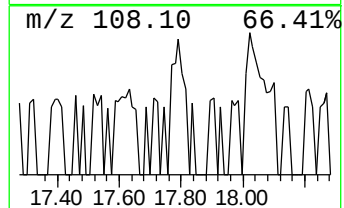
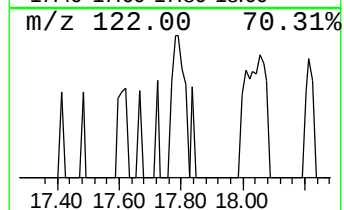
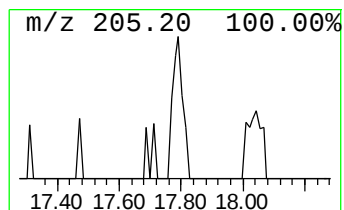
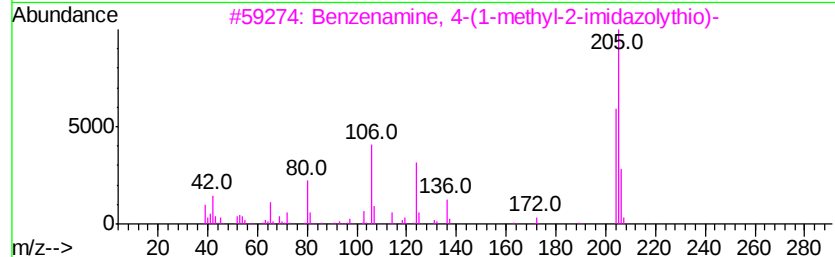
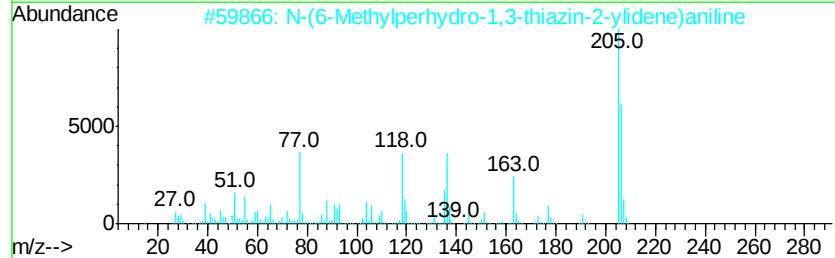
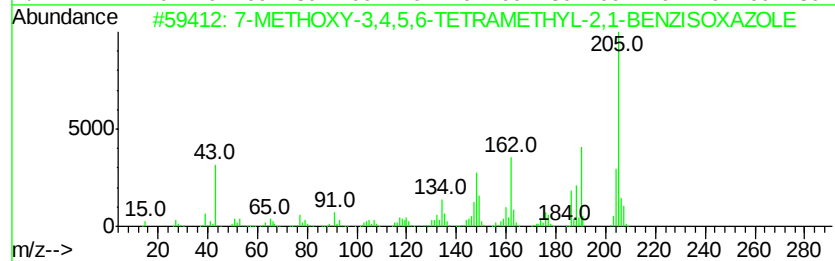
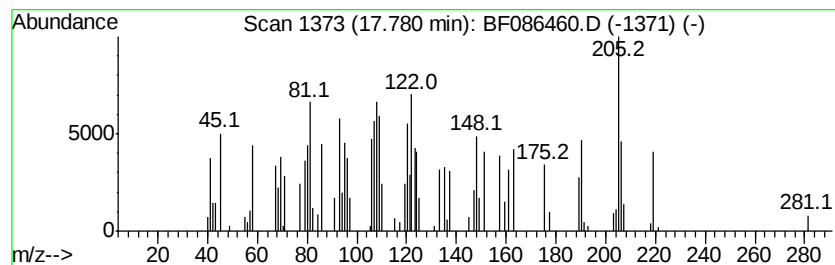
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 unknown17.78 Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.78	2.24 ng	57755	Perylene-d12	15.37

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	7-METHOXY-3,4,5,6-TETRAMETHYL-2,...	205	C12H15NO2	1000240-32-4	25
2		N-(6-Methylperhydro-1,3-thiazin-...	206	C11H14N2S	010554-33-3	14
3		Benzenamine, 4-(1-methyl-2-imida...	205	C10H11N3S	133303-54-5	14
4		.alpha.-Cedrene oxide	220	C15H24O	1000159-39-1	12
5		9-Cedranone	220	C15H24O	1000156-23-2	12



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF042716\
Data File : BF086460.D
Acq On : 28 Apr 2016 8:33
Operator : UM/SJ
Sample : H2736-04
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
RICHMOND-AMBOY-SB01(0.5-10.0)

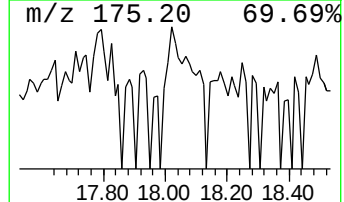
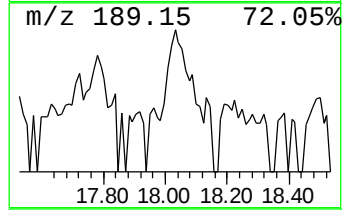
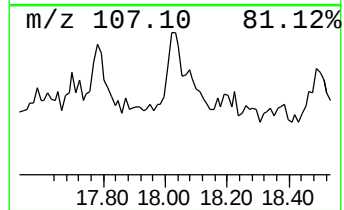
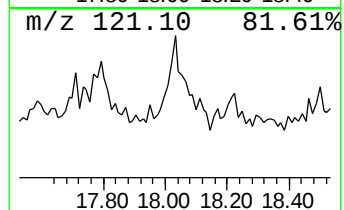
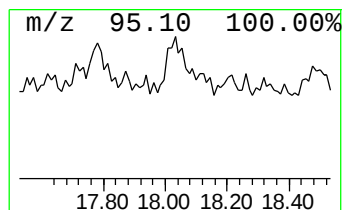
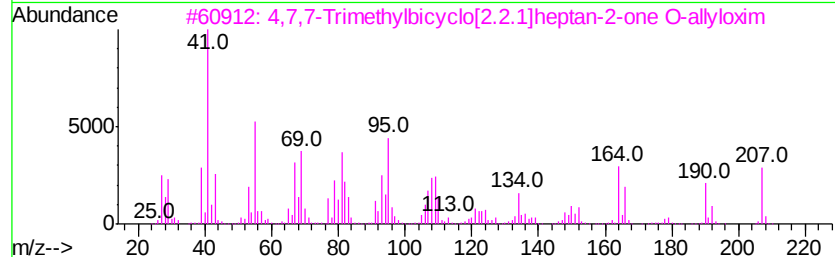
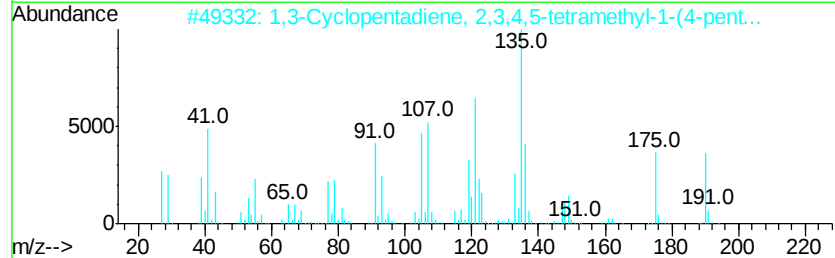
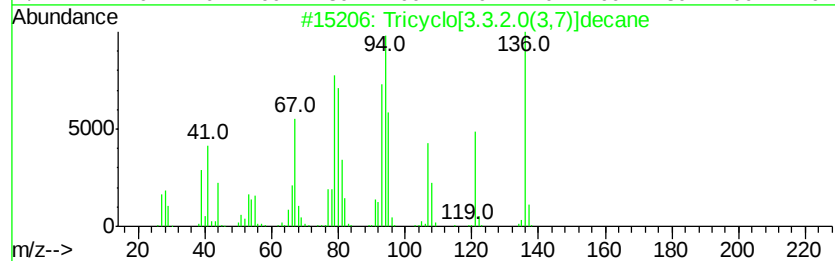
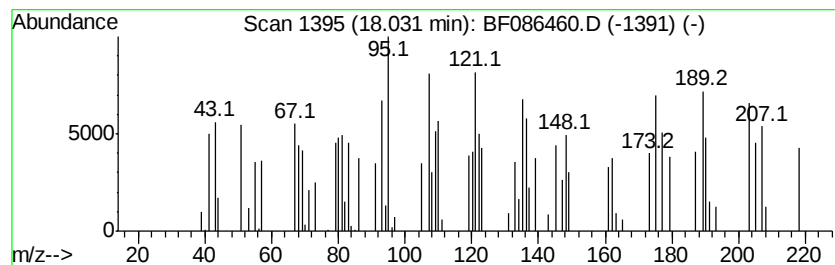
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 unknown18.03 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.03	3.08 ng	79495	Perylene-d12	15.37

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tricyclo[3.3.2.0(3,7)]decane	136	C10H16	049700-60-9	25
2		1,3-Cyclopentadiene, 2,3,4,5-tet...	190	C14H22	1000161-28-7	25
3		4,7,7-Trimethylbicyclo[2.2.1]hep...	207	C13H21NO	1000210-89-8	25
4		Tricyclo[4.4.0.0(2,8)]decane	136	C10H16	049700-59-6	22
5		Isobornyl thiocynoacetate	253	C13H19NO2S	000115-31-1	14



Data Path : Z:\HPCHEM1\BNA F\DATA\BF042716\
Data File : BF086460.D
Acq On : 28 Apr 2016 8:33
Operator : UM/SJ
Sample : H2736-04
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
RICHMOND-AMBOY-SB01(0.5-10.0)

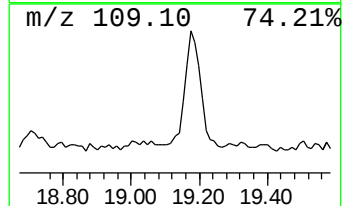
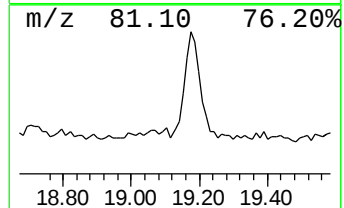
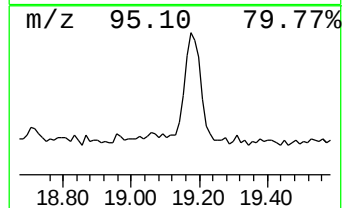
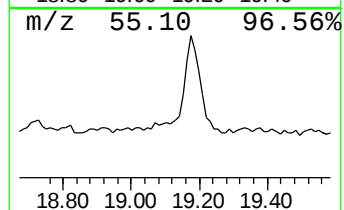
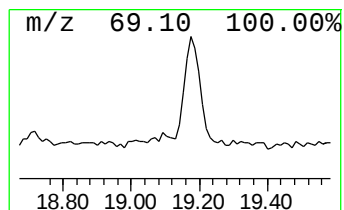
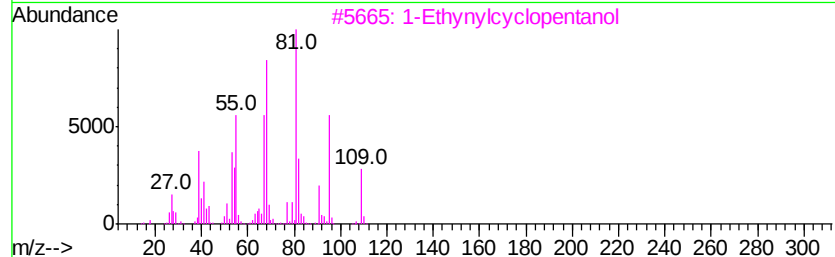
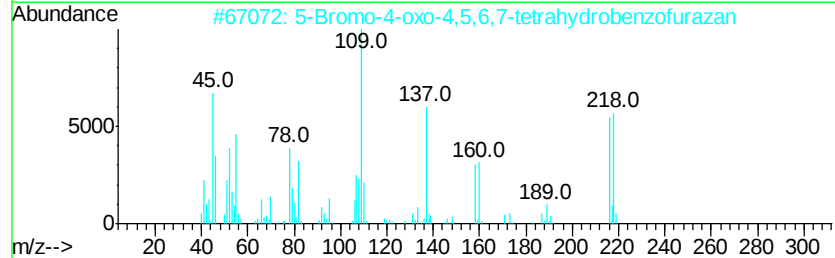
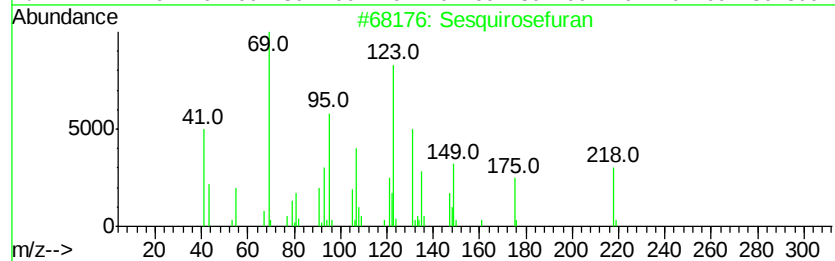
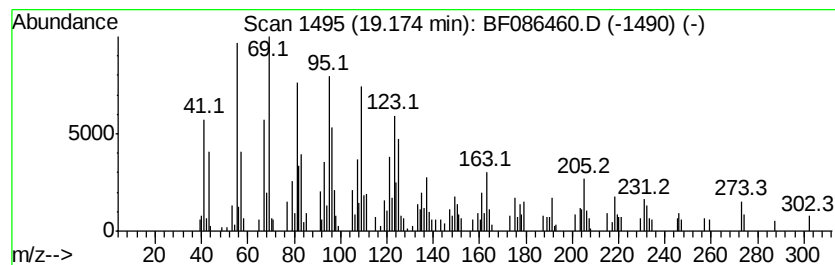
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF042716.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 Sesquirosefuran Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.17	14.37 ng	370617	Perylene-d12	15.37

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Sesquirosefuran	218	C15H22O	039007-93-7	72
2		5-Bromo-4-oxo-4,5,6,7-tetrahydro...	216	C6H5BrN2O2	300574-36-1	53
3		1-Ethynylcyclopentanol	110	C7H10O	017356-19-3	43
4		2,4a,8,8-Tetramethyldecahydrocyc...	206	C15H26	074022-04-1	42
5		2-(4a,8-Dimethyl-6-oxo-1,2,3,4,4...	234	C15H22O2	1000190-21-8	38



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF042716\
 Data File : BF086460.D
 Acq On : 28 Apr 2016 8:33
 Operator : UM/SJ
 Sample : H2736-04
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RICHMOND-AMBOY-SB01(0.5-10.0)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF042716.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...	4.99	3.1	ng	112503	1	6.81	731362	20.0
unknown6.58	6.58	109.3	ng	3996100	1	6.81	731362	20.0
Hexanoic acid, 2-...	7.47	3.5	ng	153693	2	8.09	871415	20.0
Hexadecane	10.28	2.4	ng	84873	3	9.85	705956	20.0
1-Docosene	13.81	8.7	ng	350531	5	13.95	806792	20.0
1-Hexadecanol	14.07	2.6	ng	106147	5	13.95	806792	20.0
Trichloroacetic a...	14.45	3.9	ng	157260	5	13.95	806792	20.0
Squalene	14.80	2.3	ng	59597	6	15.37	515882	20.0
Hexacosane	15.05	2.3	ng	60274	6	15.37	515882	20.0
Octacosane	15.81	2.9	ng	75507	6	15.37	515882	20.0
unknown17.47	17.47	2.6	ng	66883	6	15.37	515882	20.0
unknown17.78	17.78	2.2	ng	57755	6	15.37	515882	20.0
unknown18.03	18.03	3.1	ng	79495	6	15.37	515882	20.0
Sesquirosefuran	19.17	14.4	ng	370617	6	15.37	515882	20.0