

Data Path : Z:\HPCHEM1\BNA F\DATA\BF051016\
 Data File : BF086873.D
 Acq On : 10 May 2016 21:37
 Operator : UM/SJ
 Sample : H2916-08
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 VITALE-FILL-CSPH3-8K

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-BF050916.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	4.853	239	242	250	rBV	418934	896728	17.32%	2.032%
2	5.287	277	280	295	rBV	4400835	5040884	97.36%	11.422%
3	5.687	313	315	325	rVB	80706	116382	2.25%	0.264%
4	6.350	370	373	375	rBV	4519731	5062447	97.77%	11.471%
5	6.464	380	383	385	rBV	3761174	5177678	100.00%	11.732%
6	6.682	400	402	408	rVB	1186238	1082994	20.92%	2.454%
7	6.842	413	416	418	rBV	4266607	3982015	76.91%	9.023%
8	7.127	440	441	450	rBV	48567	129273	2.50%	0.293%
9	7.253	450	452	455	rBV	2405457	3598177	69.49%	8.153%
10	7.390	462	464	468	rVB	42670	56437	1.09%	0.128%
11	7.973	512	515	517	rBV	1274861	1236023	23.87%	2.801%
12	9.048	607	609	612	rBV	4067927	4294159	82.94%	9.730%
13	9.448	642	644	648	rBV	794629	707141	13.66%	1.602%
14	9.665	661	663	665	rBV	135748	140292	2.71%	0.318%
15	9.722	665	668	670	rBV	1161760	1177379	22.74%	2.668%
16	9.893	680	683	686	rBV2	28948	65051	1.26%	0.147%
17	10.156	705	706	708	rVB	234762	187862	3.63%	0.426%
18	10.373	722	725	729	rBV	72475	131703	2.54%	0.298%
19	10.511	734	737	739	rBV	2479414	2662333	51.42%	6.033%
20	10.625	746	747	752	rBV	201179	271225	5.24%	0.615%
21	11.071	785	786	792	rVB	97626	157397	3.04%	0.357%
22	11.196	795	797	800	rBV	1114738	964866	18.64%	2.186%
23	11.493	821	823	826	rVB	38463	52372	1.01%	0.119%
24	11.745	843	845	848	rBV	115430	166589	3.22%	0.377%
25	12.522	911	913	919	rVB	71040	83580	1.61%	0.189%
26	12.785	933	936	938	rBV	2995316	3586209	69.26%	8.126%
27	13.014	954	956	962	rBV3	20653	62773	1.21%	0.142%
28	13.688	1013	1015	1023	rBV	367806	779145	15.05%	1.765%
29	13.825	1025	1027	1029	rBV	771192	809505	15.63%	1.834%
30	14.271	1063	1066	1068	rBV	114799	178451	3.45%	0.404%
31	14.339	1070	1072	1076	rVB	32604	55858	1.08%	0.127%
32	15.197	1144	1147	1153	rVB	564638	709423	13.70%	1.607%
33	15.700	1187	1191	1198	rBV4	39256	143244	2.77%	0.325%
34	15.974	1213	1215	1218	rVV	28657	52999	1.02%	0.120%

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ClientSampleId :
VITALE-FILL-CSPH3-8K

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-BF050916.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	16.043	1218	1221	1228	rVB2	30881	73370	1.42%	0.166%
36	16.580	1263	1268	1273	rVB3	41350	139778	2.70%	0.317%
37	17.311	1329	1332	1337	rVB	32488	100885	1.95%	0.229%

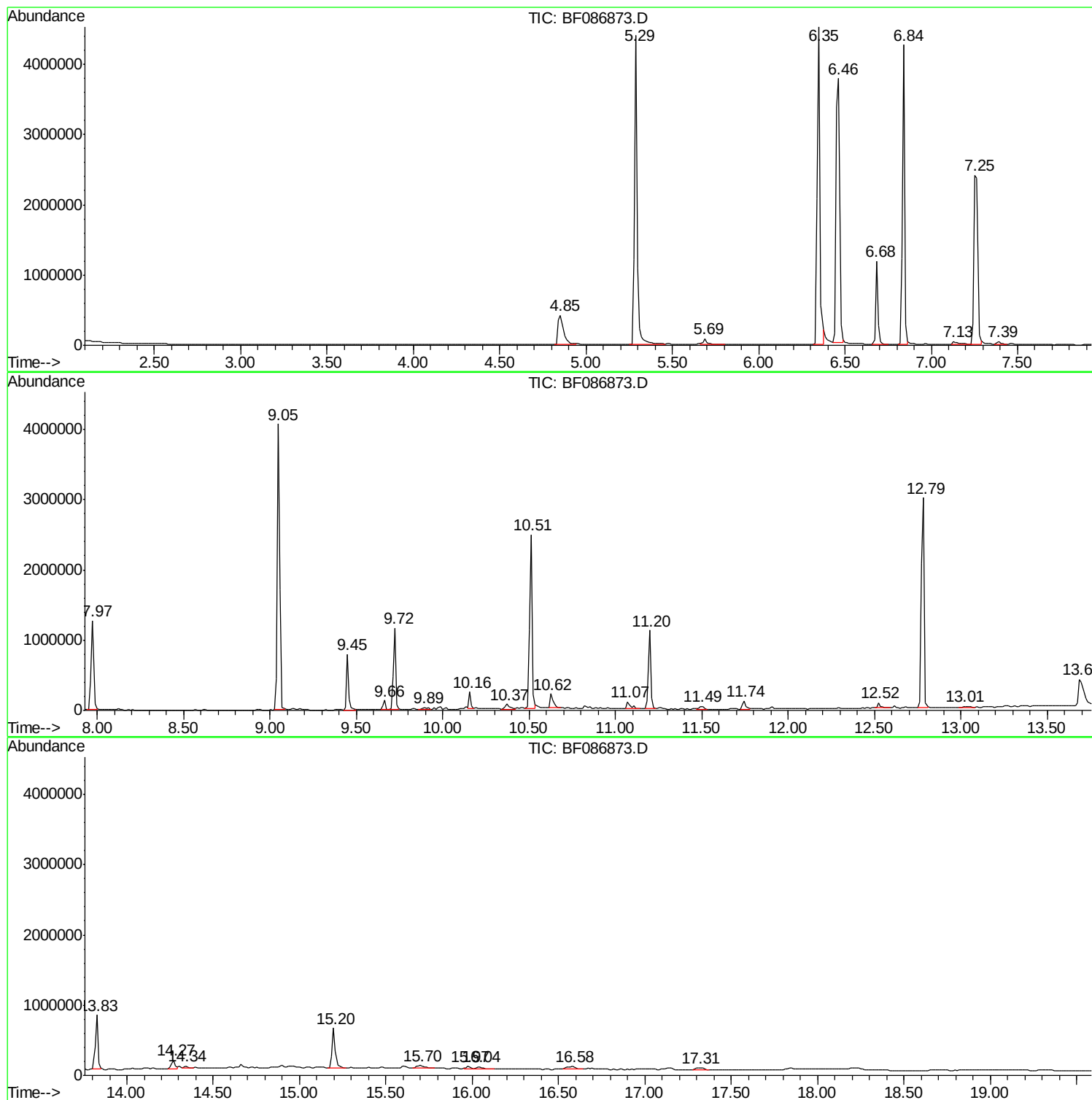
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF050916.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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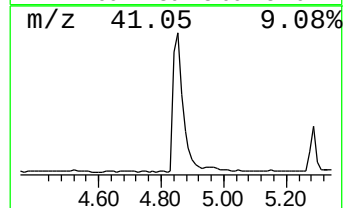
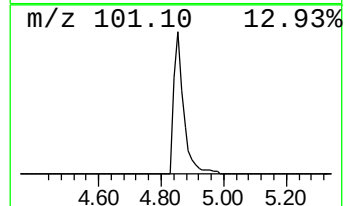
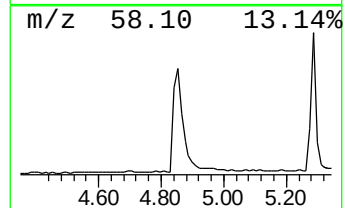
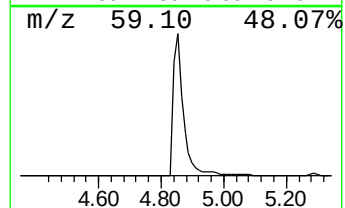
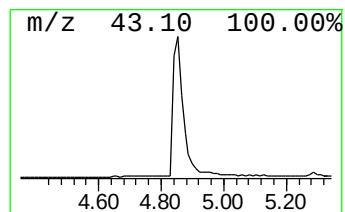
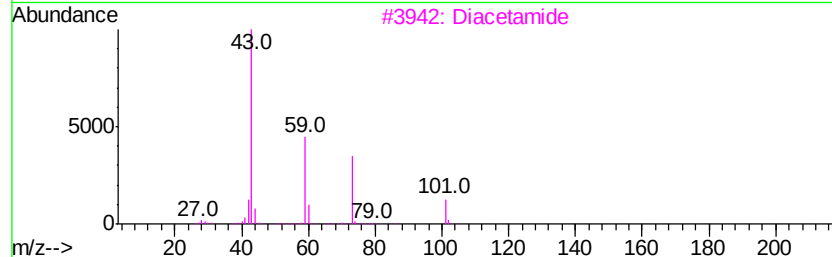
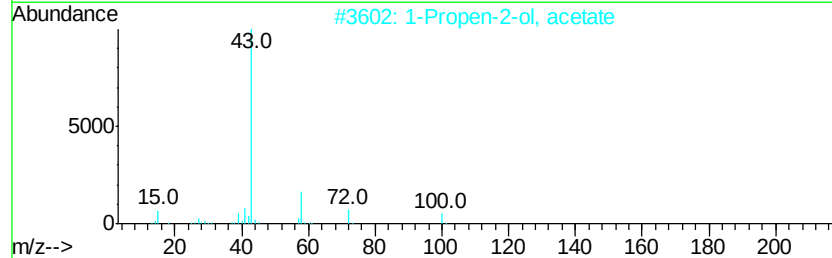
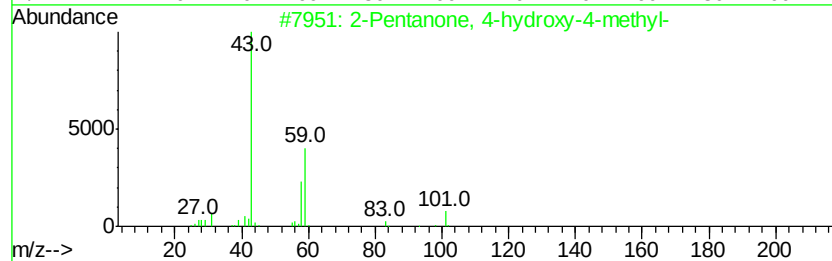
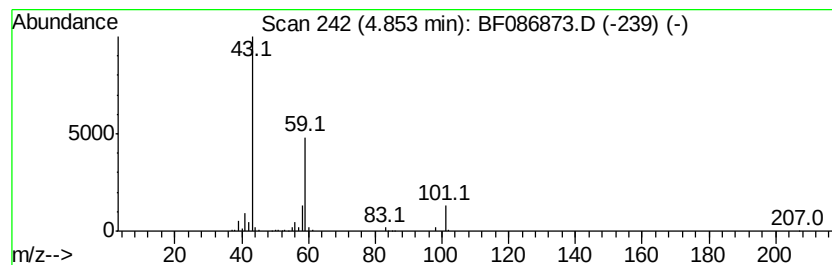
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.85	16.56 ng	896728	1,4-Dichlorobenzene-d4	6.68

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2		1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10
3		Diacetamide	101	C4H7NO2	000625-77-4	9
4		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
5		Acetic acid, 2-propenyl ester	100	C5H8O2	000591-87-7	9



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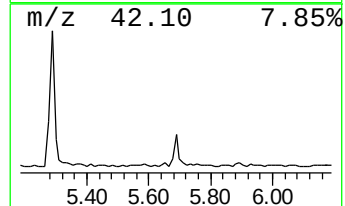
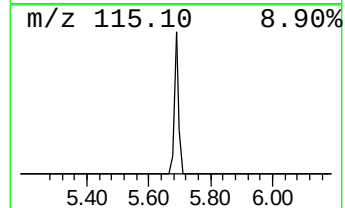
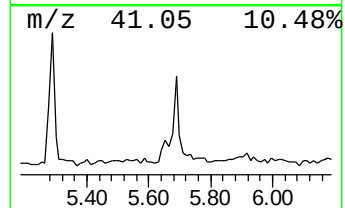
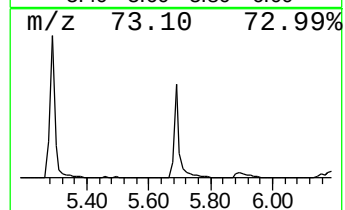
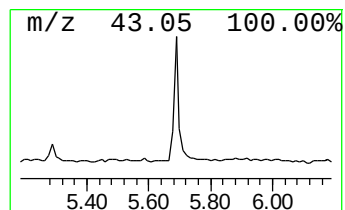
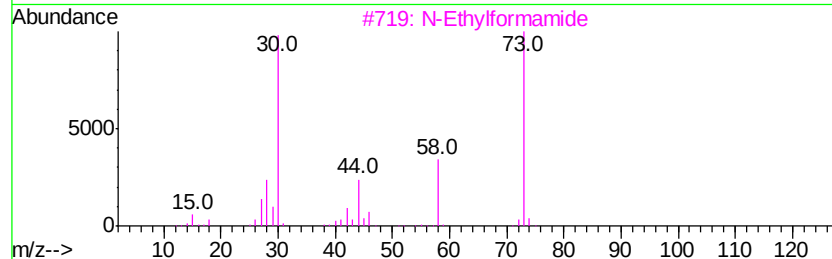
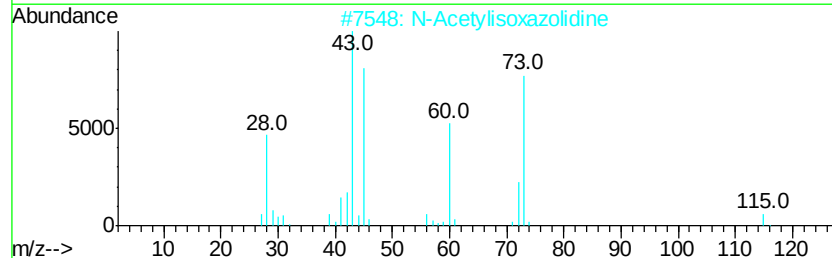
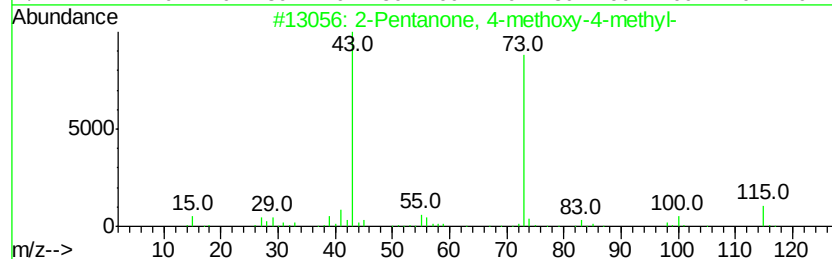
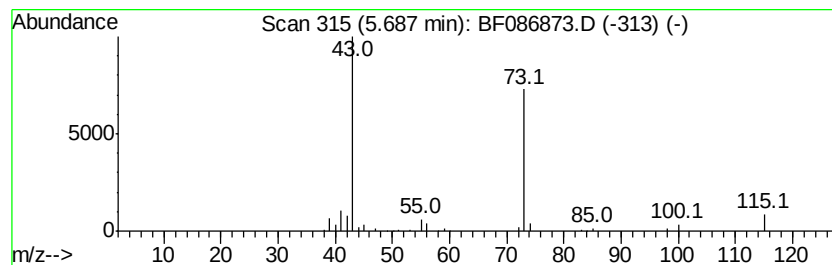
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 2-Pentanone, 4-methoxy-4-me... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.69	2.15 ng	116382	1,4-Dichlorobenzene-d4	6.68

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-methoxy-4-methyl-	130	C7H14O2	000107-70-0	83
2		N-Acetylisoaxazolidine	115	C5H9N2O2	115615-36-6	33
3		N-Ethylformamide	73	C3H7NO	000627-45-2	12
4		Borane, dimethoxy-	74	C2H7BO2	004542-61-4	9
5		Acetamide, N-methyl-	73	C3H7NO	000079-16-3	9



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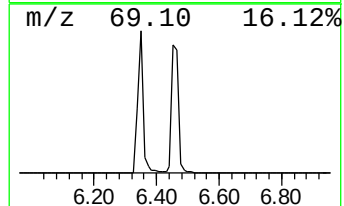
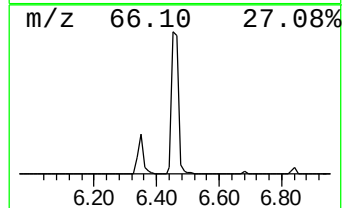
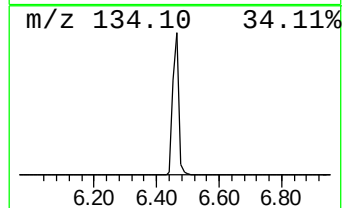
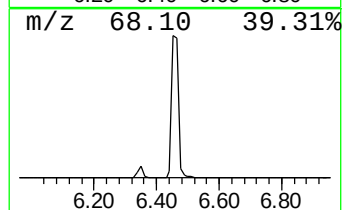
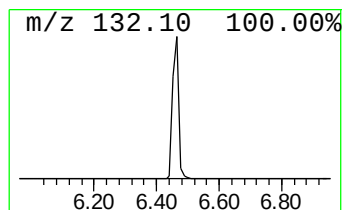
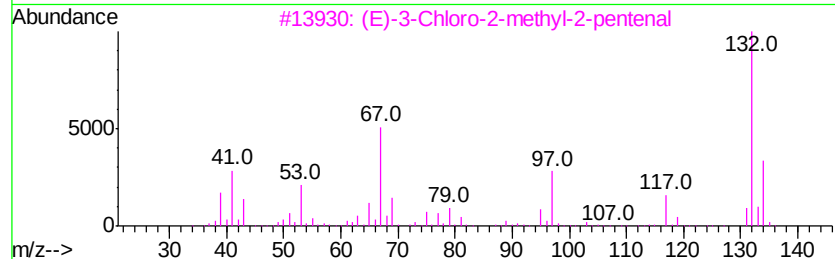
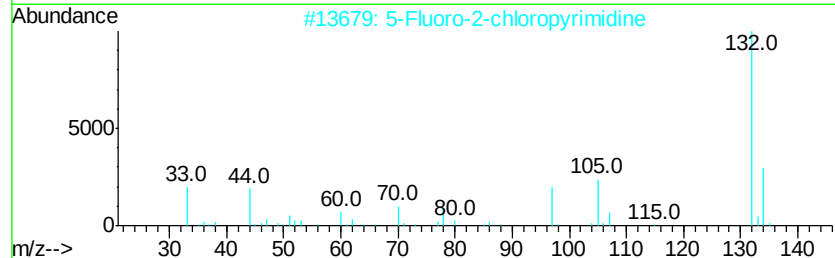
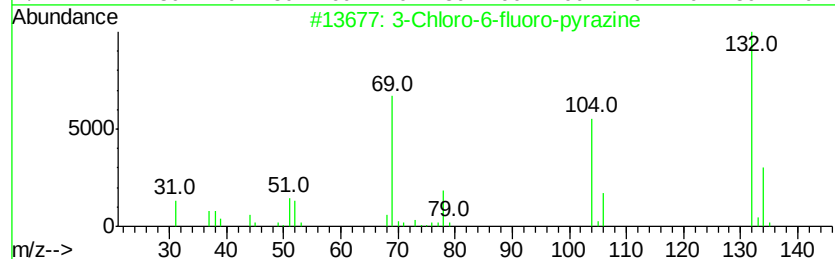
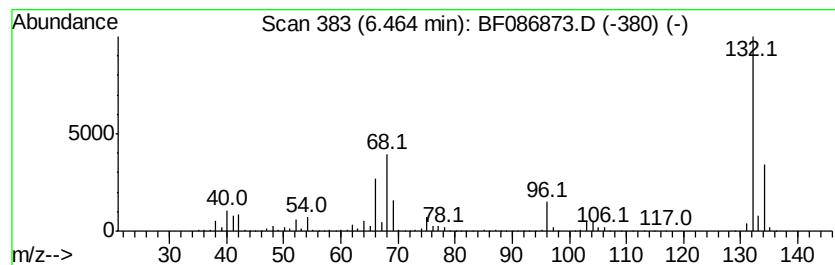
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 unknown6.46 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.46	95.62 ng	5177680	1,4-Dichlorobenzene-d4	6.68

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
3		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	12
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
5		Benzene, 1,2,4-trifluoro-	132	C6H3F3	000367-23-7	9



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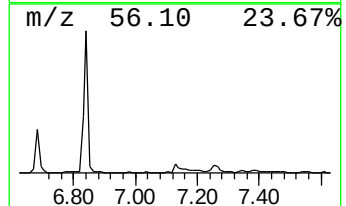
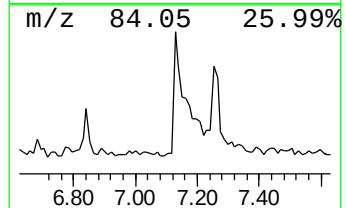
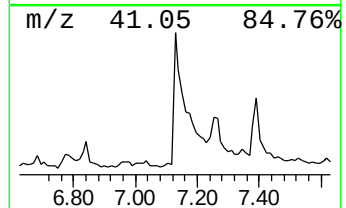
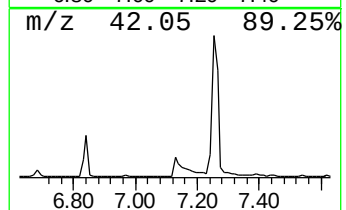
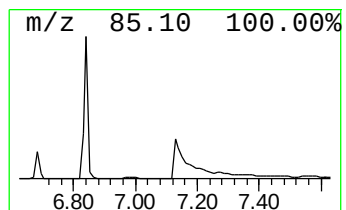
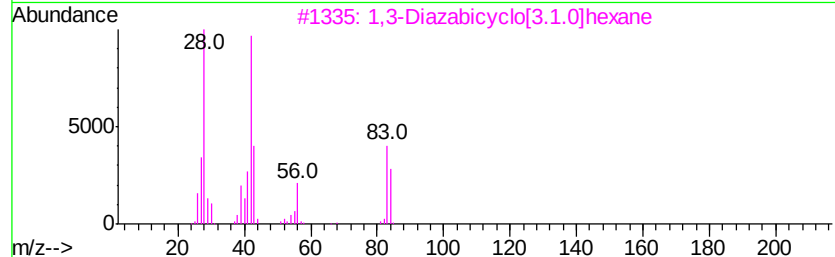
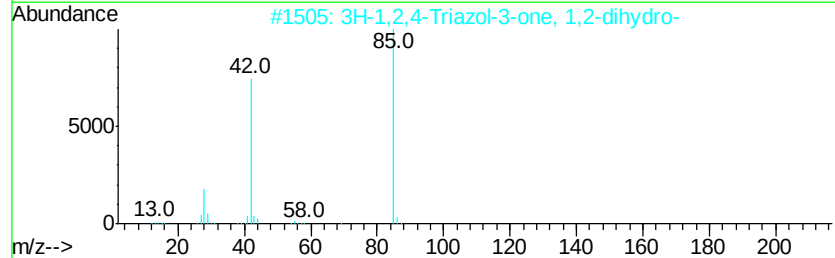
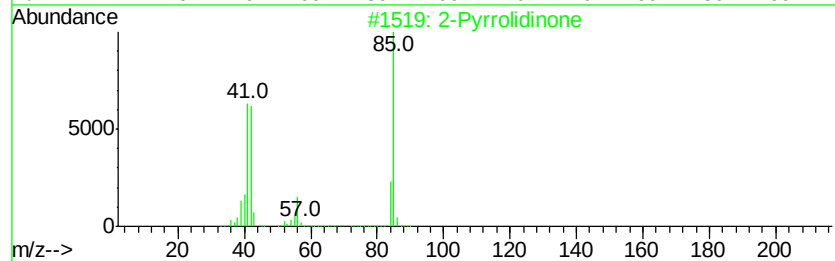
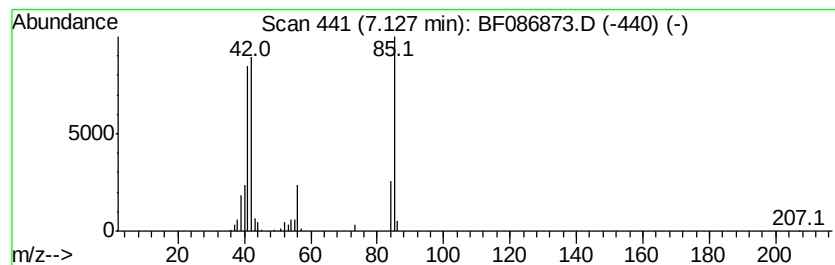
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 2-Pyrrolidinone Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.13	2.39 ng	129273	1,4-Dichlorobenzene-d4	6.68

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pyrrolidinone	85	C4H7NO	000616-45-5	91
2		3H-1,2,4-Triazol-3-one, 1,2-dihy...	85	C2H3N3O	000930-33-6	53
3		1,3-Diazabicyclo[3.1.0]hexane	84	C4H8N2	017038-28-7	43
4		1-Hexene	84	C6H12	000592-41-6	38
5		2-Oxetanone, 4-methylene-	84	C4H4O2	000674-82-8	25



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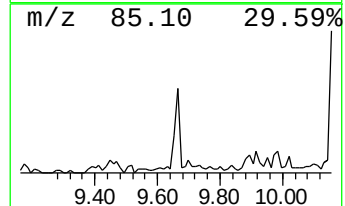
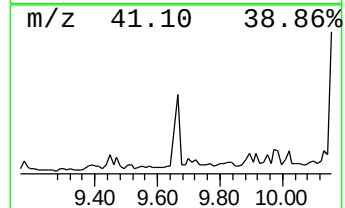
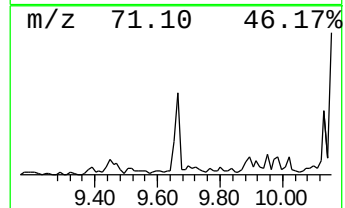
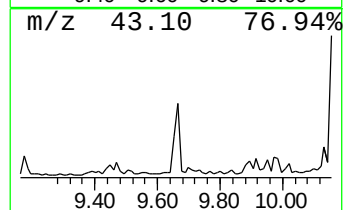
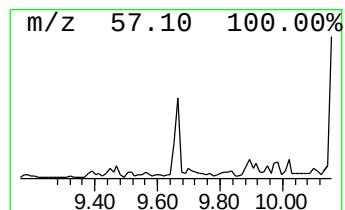
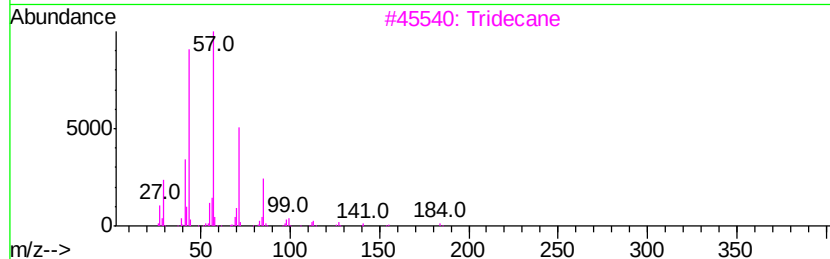
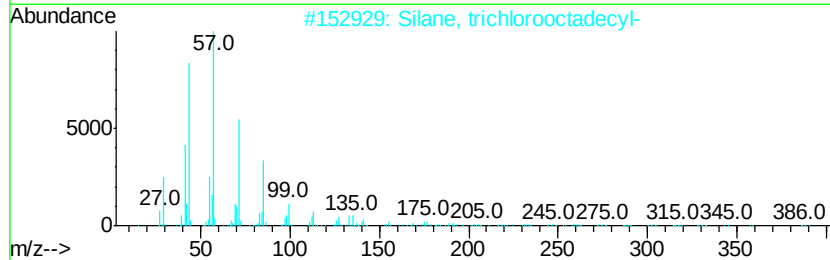
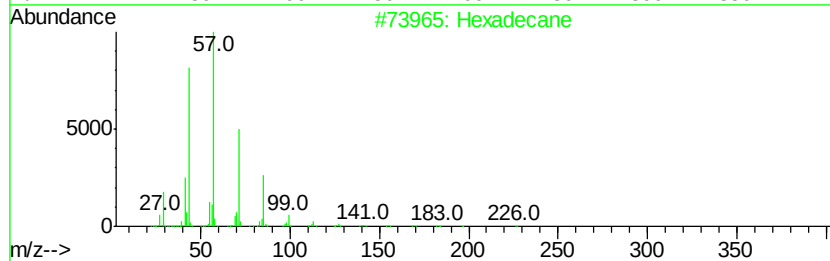
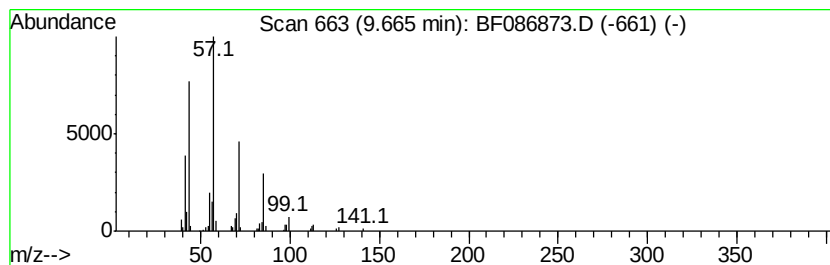
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Hexadecane Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.66	2.38 ng	140292	Acenaphthene-d10	9.72

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecane	226	C16H34	000544-76-3	90
2		Silane, trichlorooctadecyl-	386	C18H37Cl3Si	000112-04-9	86
3		Tridecane	184	C13H28	000629-50-5	86
4		Tetradecane	198	C14H30	000629-59-4	80
5		Eicosane	282	C20H42	000112-95-8	64



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Operator : UM/SJ
Sample : H2916-08
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
VITALE-FILL-CSPH3-8K

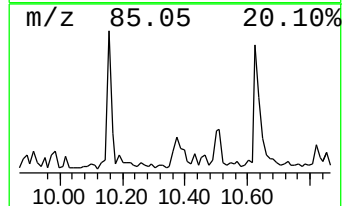
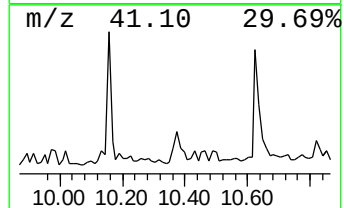
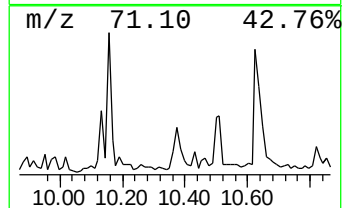
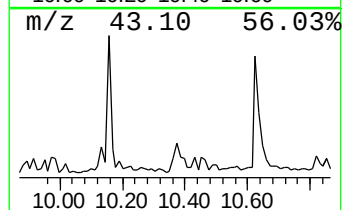
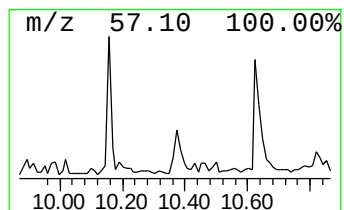
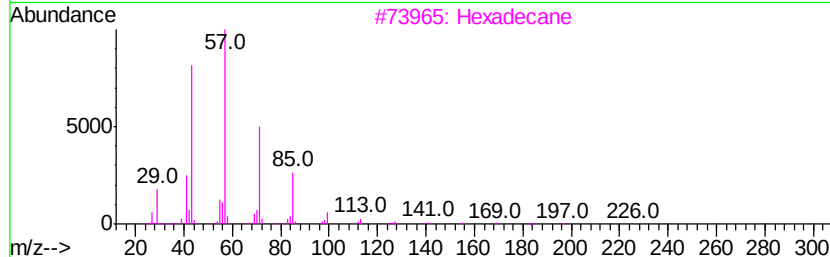
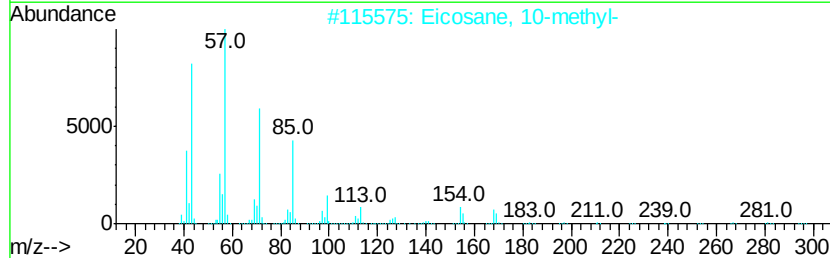
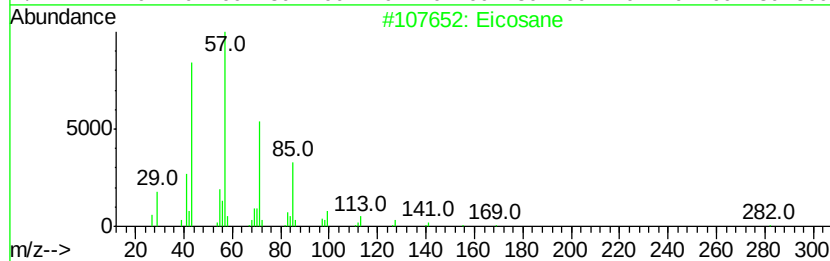
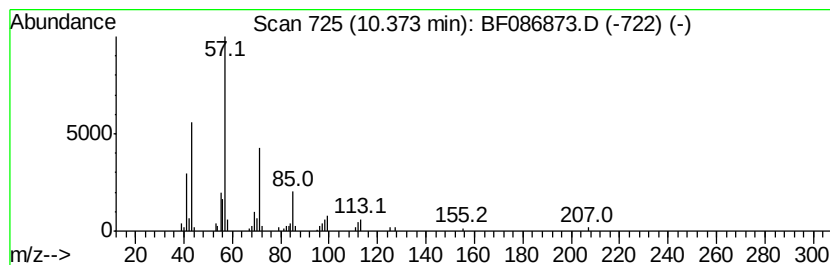
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Eicosane Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.37	2.24 ng	131703	Acenaphthene-d10	9.72

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	80
2		Eicosane, 10-methyl-	296	C21H44	054833-23-7	72
3		Hexadecane	226	C16H34	000544-76-3	72
4		Tridecane, 6-methyl-	198	C14H30	013287-21-3	72
5		Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	64



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF051016\
Data File : BF086873.D
Acq On : 10 May 2016 21:37
Operator : UM/SJ
Sample : H2916-08
Misc :
ALS Vial : 22 Sample Multiplier: 1

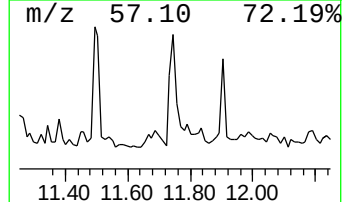
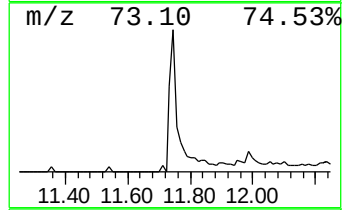
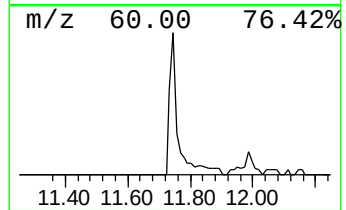
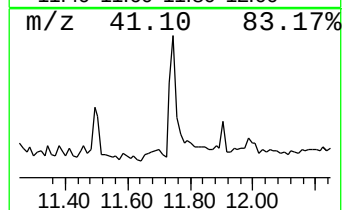
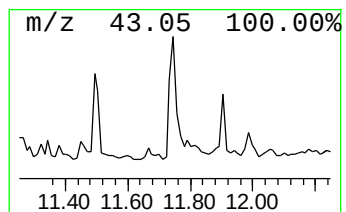
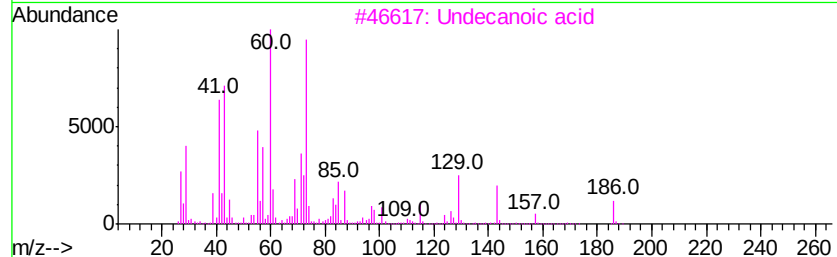
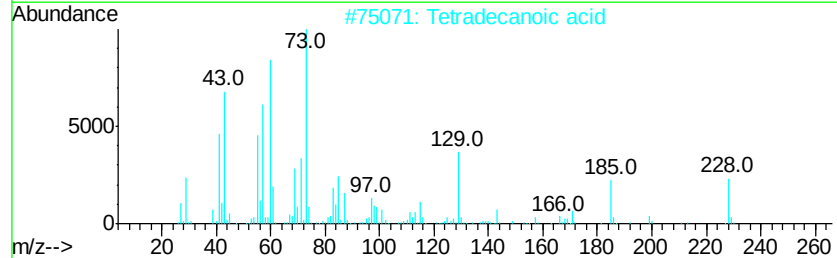
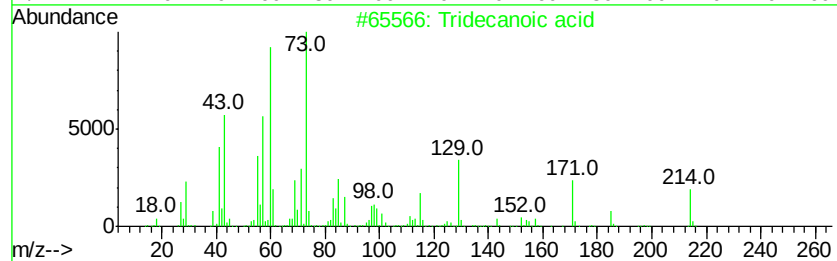
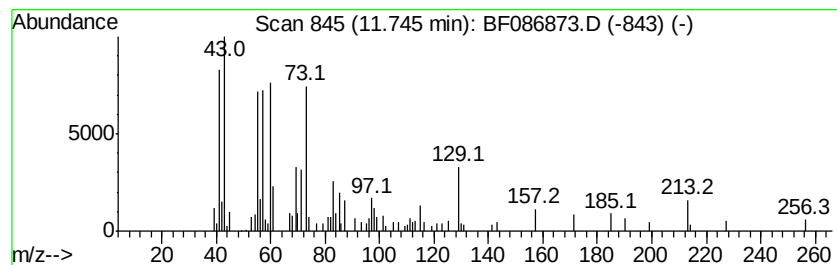
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ClientSampleId :
VITALE-FILL-CSPH3-8K

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF050916.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Tridecanoic acid Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD		R.T.
11.75	3.45 ng	166589	Phenanthrene-d10		11.20
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tridecanoic acid	214	C13H26O2	000638-53-9	80
2	Tetradecanoic acid	228	C14H28O2	000544-63-8	64
3	Undecanoic acid	186	C11H22O2	000112-37-8	59
4	Dodecanoic acid	200	C12H24O2	000143-07-7	59
5	Pentadecanoic acid	242	C15H30O2	001002-84-2	58



Data Path : Z:\HPCHEM1\BNA F\DATA\BF051016\
Data File : BF086873.D
Acq On : 10 May 2016 21:37
Operator : UM/SJ
Sample : H2916-08
Misc :
ALS Vial : 22 Sample Multiplier: 1

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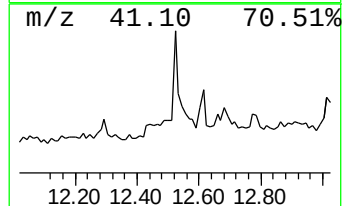
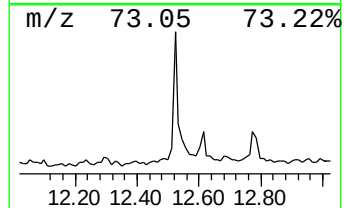
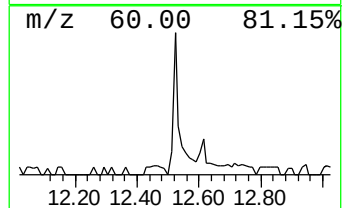
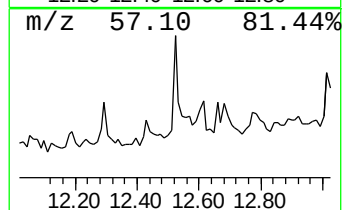
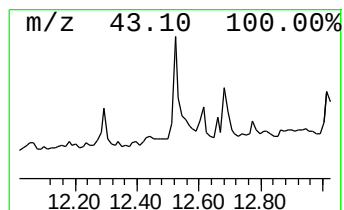
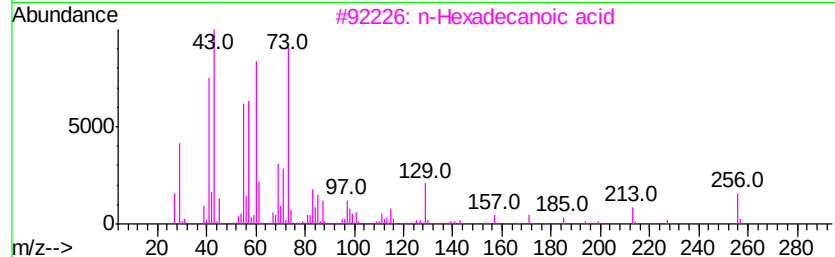
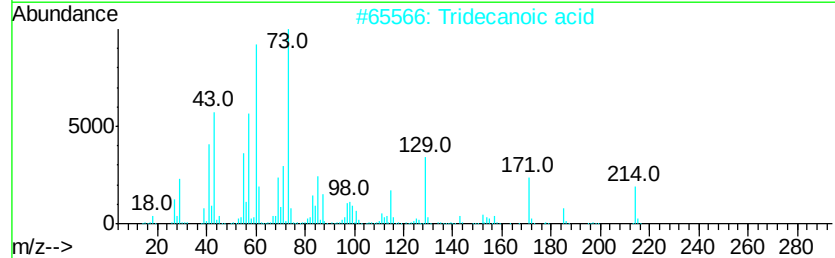
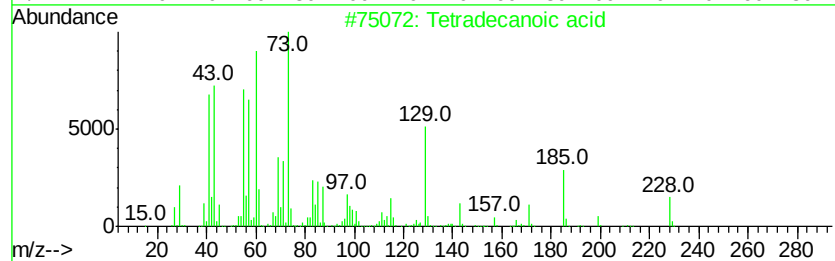
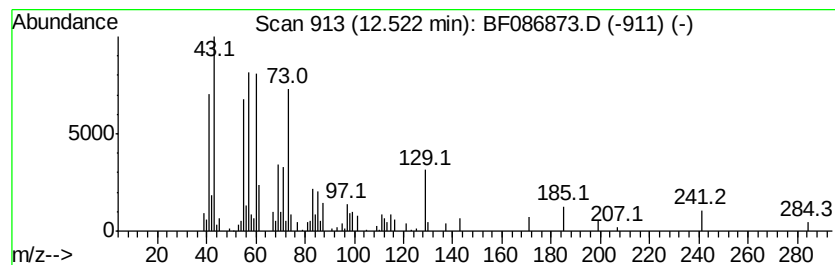
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Tetradecanoic acid Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.52	2.06 ng	83580	Chrysene-d12	13.83

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetradecanoic acid	228	C14H28O2	000544-63-8	74
2			Tridecanoic acid	214	C13H26O2	000638-53-9	72
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	72
4			Pentadecanoic acid	242	C15H30O2	001002-84-2	72
5			Undecanoic acid	186	C11H22O2	000112-37-8	53



Data Path : Z:\HPCHEM1\BNA F\DATA\BF051016\
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Acq On : 10 May 2016 21:37
Operator : UM/SJ
Sample : H2916-08
Misc :
ALS Vial : 22 Sample Multiplier: 1

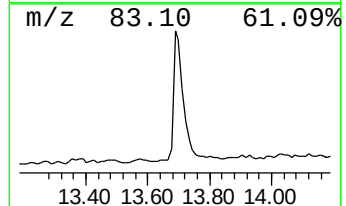
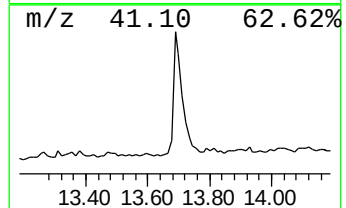
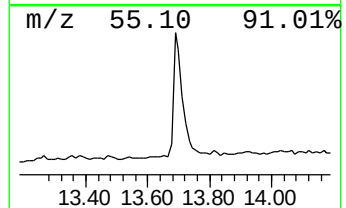
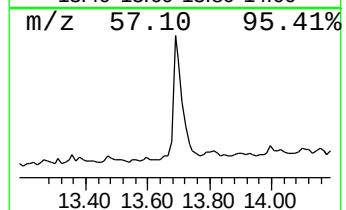
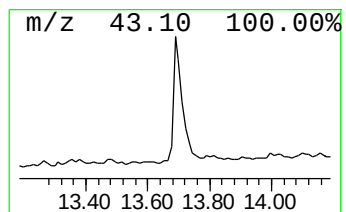
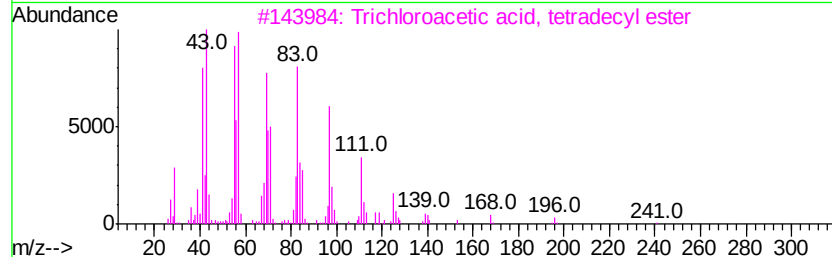
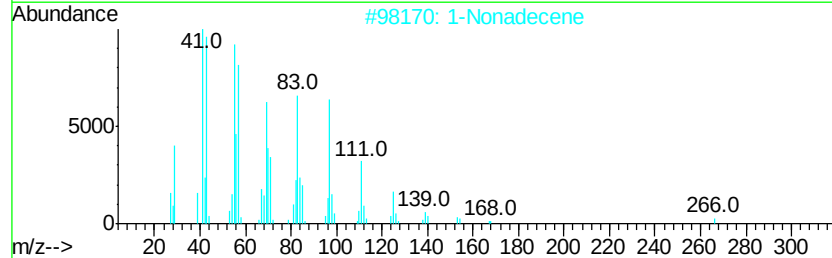
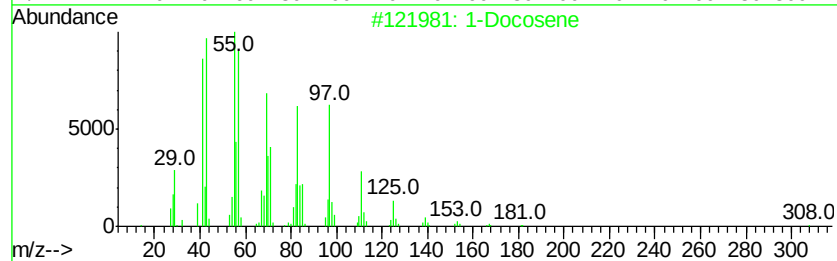
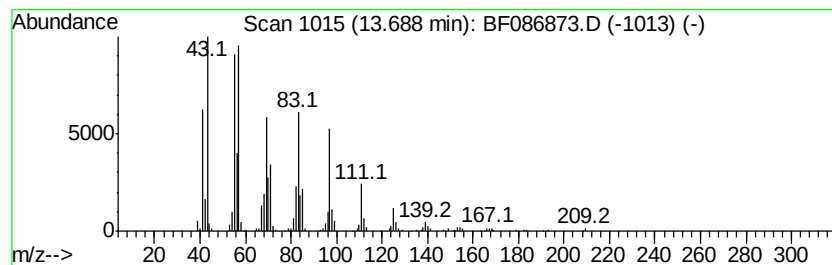
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 1-Docosene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.69	19.25 ng	779145	Chrysene-d12	13.83	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Docosene	308	C22H44	001599-67-3	94
2	1-Nonadecene	266	C19H38	018435-45-5	90
3	Trichloroacetic acid, tetradecyl...	358	C16H29Cl3O2	074339-52-9	87
4	Pentafluoropropionic acid, hexad...	388	C19H33F5O2	006222-07-7	83
5	Pentafluoropropionic acid, octad...	416	C21H37F5O2	1000280-07-7	83



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF051016\
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Sample : H2916-08
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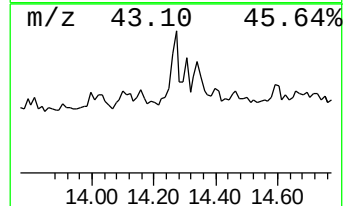
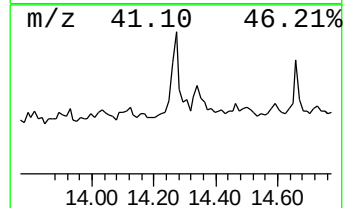
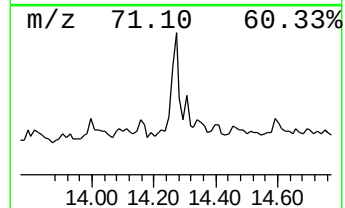
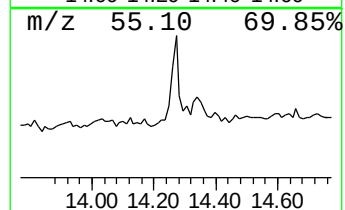
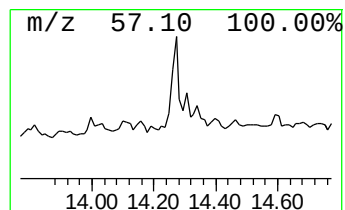
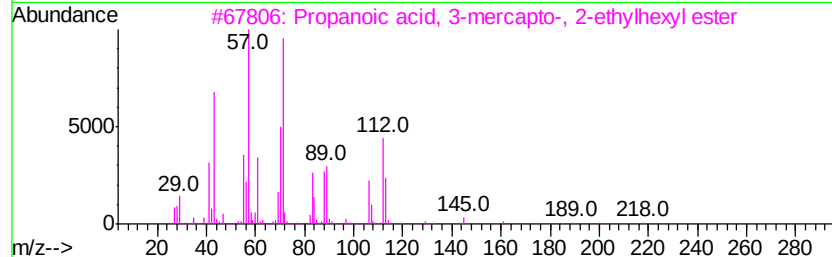
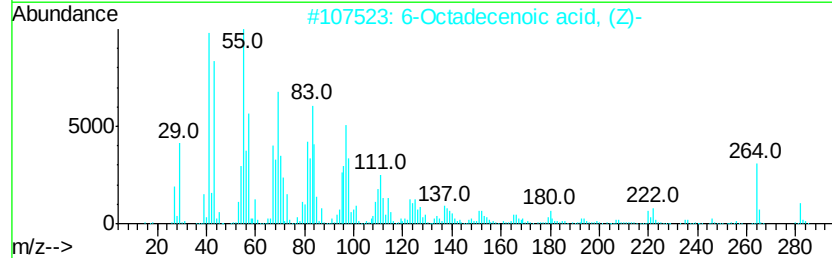
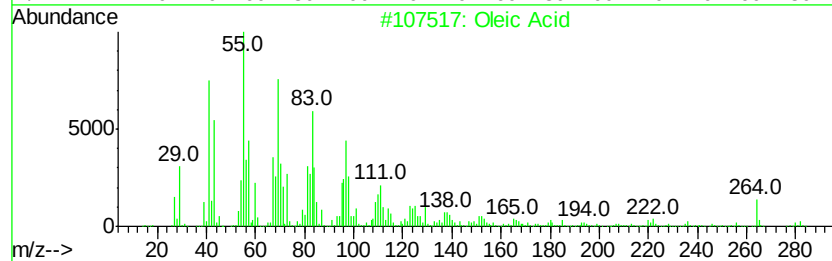
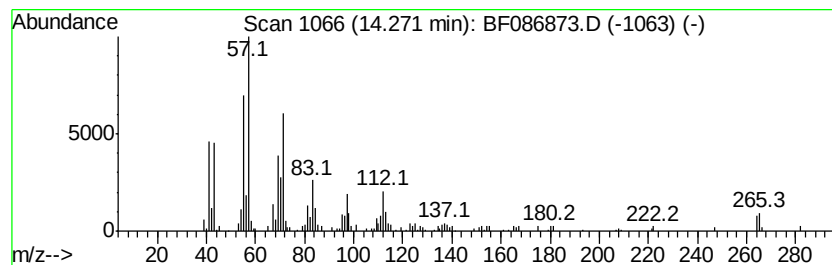
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Oleic Acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.27	4.41 ng	178451	Chrysene-d12	13.83	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Oleic Acid	282	C18H34O2	000112-80-1	94
2	6-Octadecenoic acid, (Z)-	282	C18H34O2	000593-39-5	80
3	Propanoic acid, 3-mercapto-, 2-e...	218	C11H22O2S	050448-95-8	47
4	9-Octadecenoic acid, (E)-	282	C18H34O2	000112-79-8	35
5	Dichloroacetic acid, heptadecyl ...	366	C19H36Cl2O2	1000282-98-2	35



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF051016\
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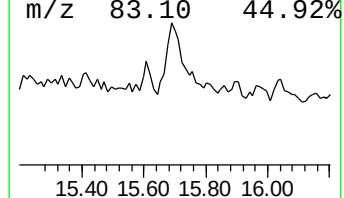
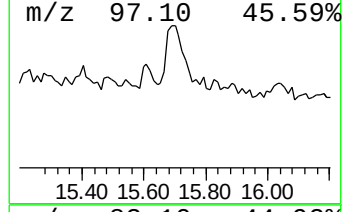
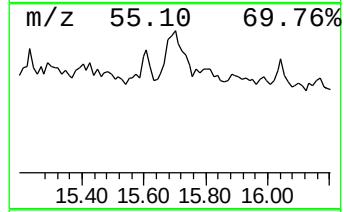
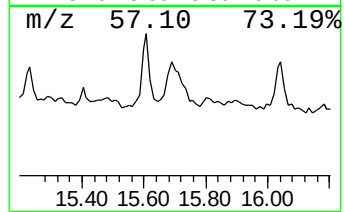
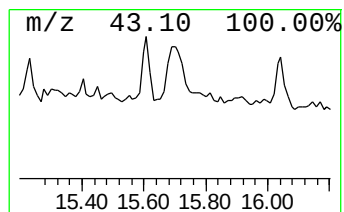
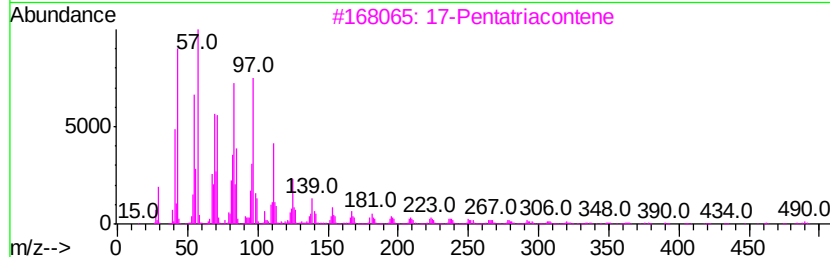
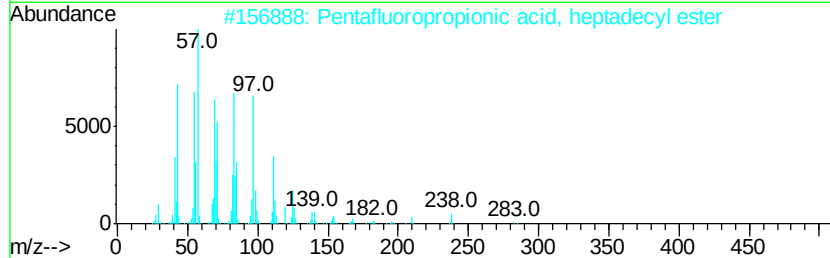
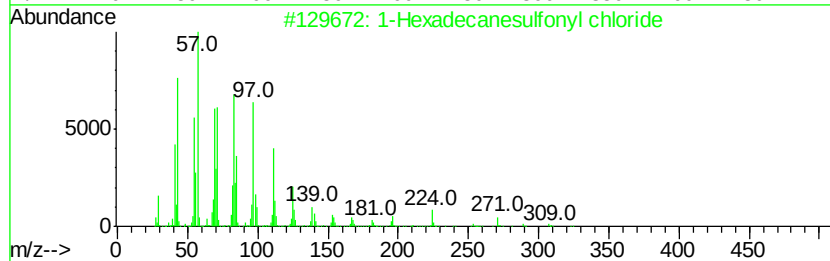
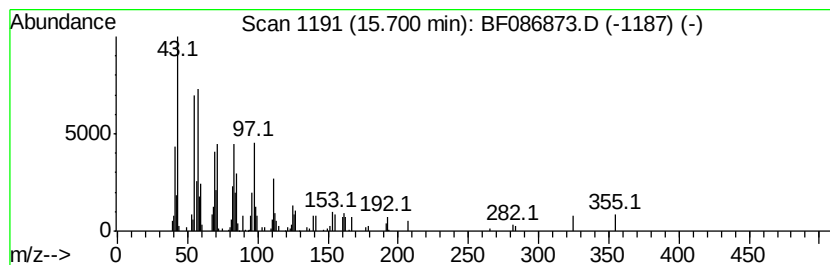
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 1-Hexadecanesulfonyl chloride Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.		
15.70	4.04 ng	143244	Perylene-d12	15.20		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Hexadecanesulfonyl chloride	324	C16H33ClO2S	038775-38-1	81	
2	Pentafluoropropionic acid, hepta...	402	C20H35F5O2	1000283-04-2	81	
3	17-Pentatriacontene	491	C35H70	006971-40-0	81	
4	Oxirane, [(dodecyloxy)methyl]-	242	C15H30O2	002461-18-9	74	
5	Heptafluorobutyric acid, n-octad...	466	C22H37F7O2	000400-57-7	74	



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF051016\
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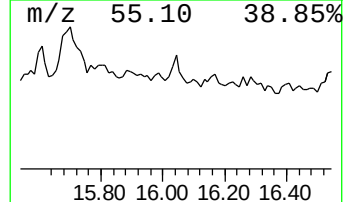
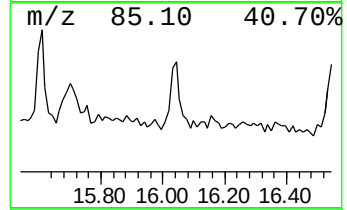
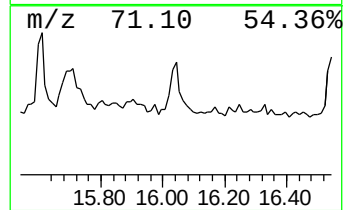
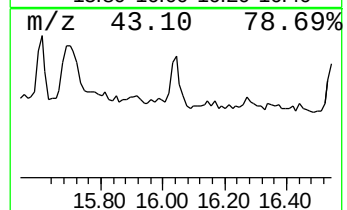
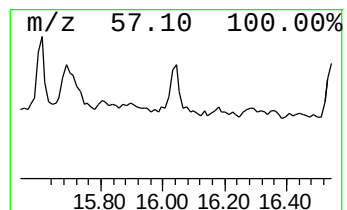
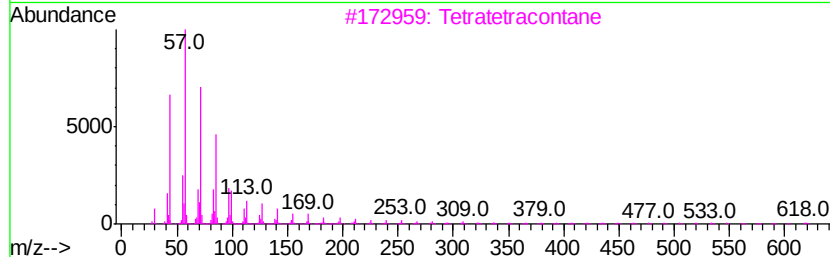
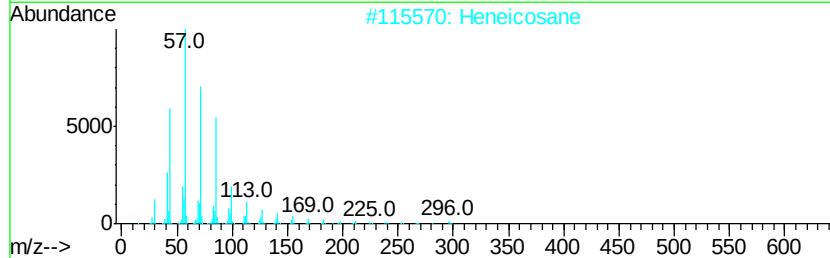
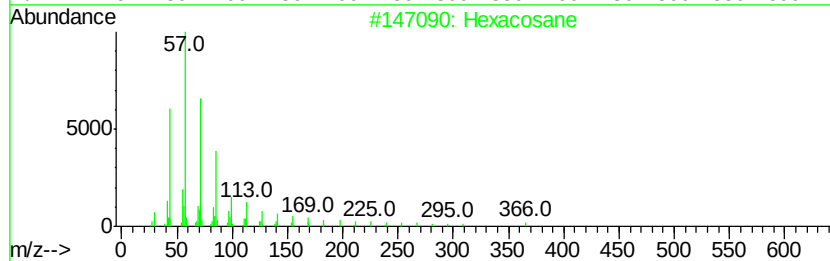
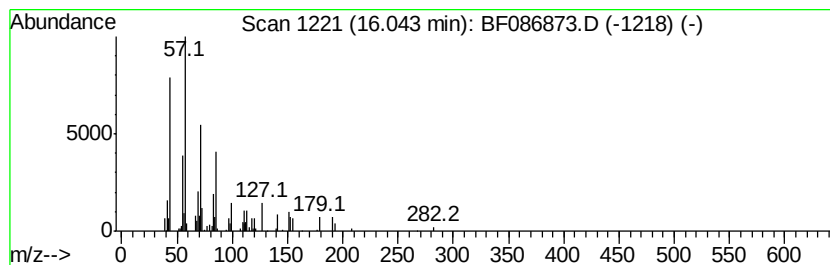
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 16 Hexacosane Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.04	2.07 ng	73370	Perylene-d12	15.20
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Hexacosane	366 C26H54	000630-01-3	72
2	Heneicosane	296 C21H44	000629-94-7	72
3	Tetratetracontane	619 C44H90	007098-22-8	72
4	Pentacosane	352 C25H52	000629-99-2	72
5	Eicosane	282 C20H42	000112-95-8	70



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF051016\
Data File : BF086873.D
Acq On : 10 May 2016 21:37
Operator : UM/SJ
Sample : H2916-08
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
VITALE-FILL-CSPH3-8K

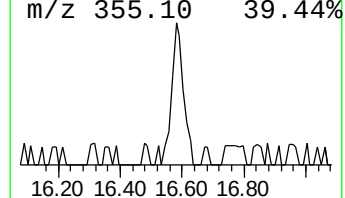
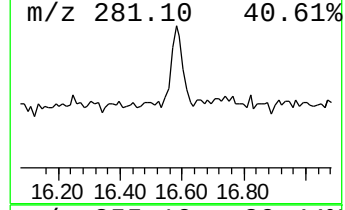
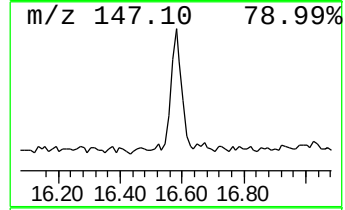
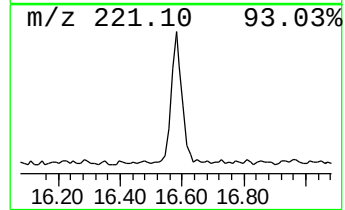
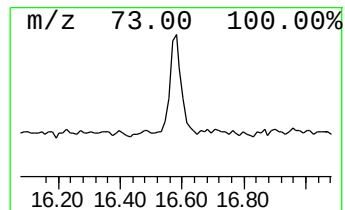
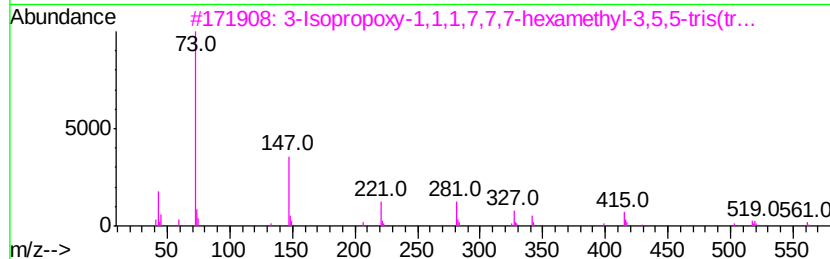
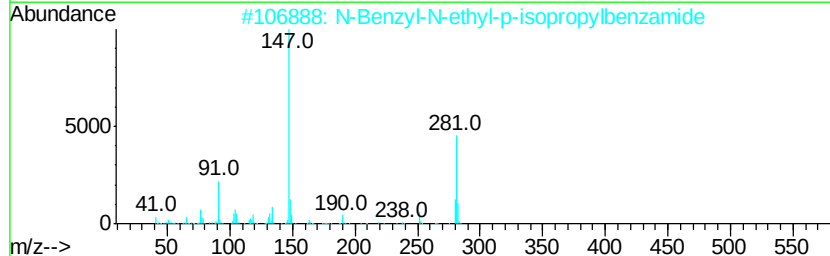
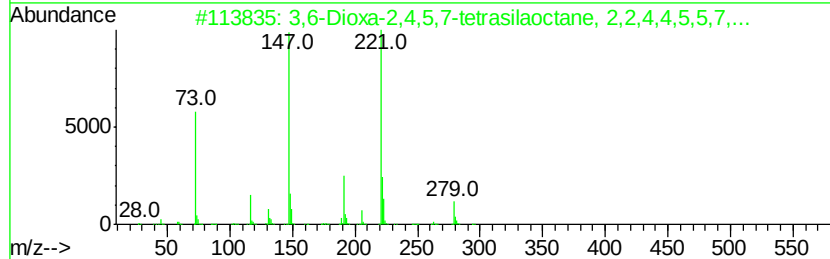
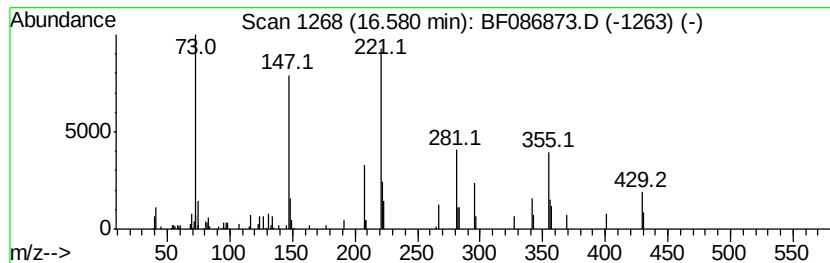
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 unknown16.58 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.58	3.94 ng	139778	Perylene-d12	15.20

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3,6-Dioxa-2,4,5,7-tetrasilaoctan...	294	C10H30O2Si4	004342-25-0	38
2			N-Benzyl-N-ethyl-p-isopropylbenz...	281	C19H23NO	015089-22-2	25
3			3-Isopropoxy-1,1,1,7,7,7-hexamet...	576	C18H52O7Si7	071579-69-6	25
4			2,6-Dimethyl-3,4-bis(trimethylsi...	311	C15H29NO2Si2	1000079-52-1	12
5			Piperidine, 1-(5-trifluoromethyl...	295	C15H16F3N3	1000268-74-7	10



Data Path : Z:\HPCHEM1\BNA F\DATA\BF051016\
Data File : BF086873.D
Acq On : 10 May 2016 21:37
Operator : UM/SJ
Sample : H2916-08
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
VITALE-FILL-CSPH3-8K

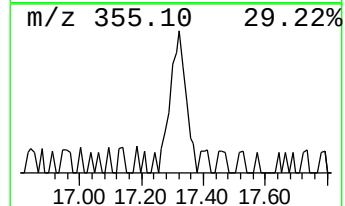
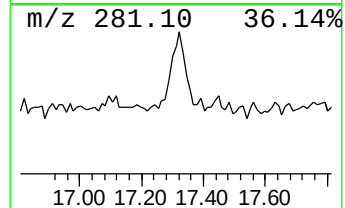
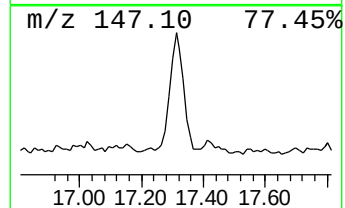
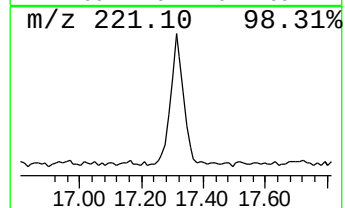
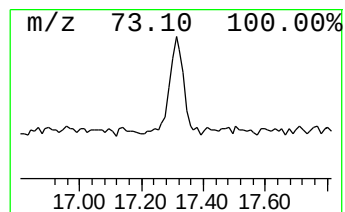
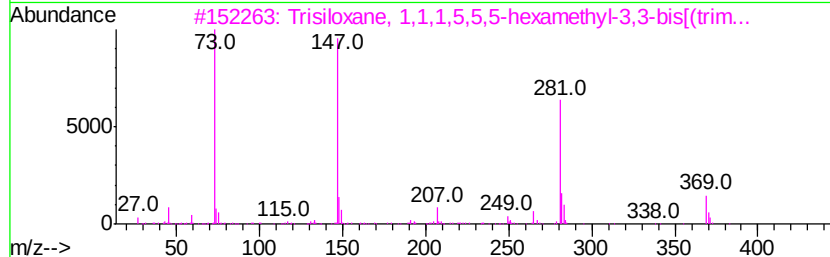
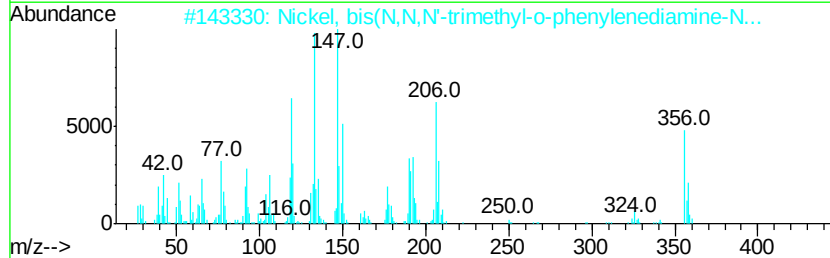
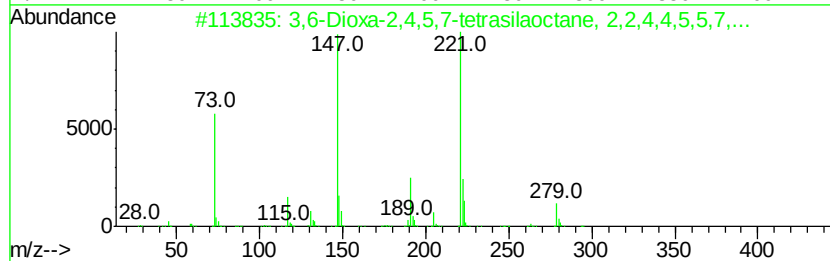
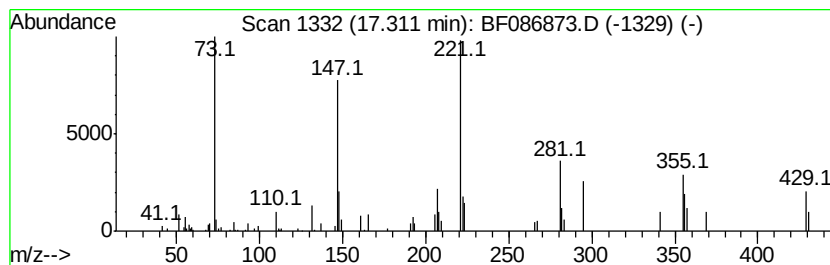
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 unknown17.31 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.31	2.84 ng	100885	Perylene-d12	15.20

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3,6-Dioxa-2,4,5,7-tetrasilaoctan...	294	C10H30O2Si4	004342-25-0	37
2			Nickel, bis(N,N,N'-trimethyl-o-p...	356	C18H26N4Ni	1000153-51-1	25
3			Trisiloxane, 1,1,1,5,5,5-hexamet...	384	C12H36O4Si5	003555-47-3	16
4			Heptasiloxane, 1,1,3,3,5,5,7,7,9...	504	C14H44O6Si7	019095-23-9	14
5			1H-1,2,3-Triazole, 4,5-diphenyl-	221	C14H11N3	005533-73-3	11



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF051016\
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 Sample : H2916-08
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 VITALE-FILL-CSPH3-8K

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF050916.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...	4.85	16.6	ng	896728	1	6.68	1082990	20.0
2-Pentanone, 4-me...	5.69	2.1	ng	116382	1	6.68	1082990	20.0
unknown6.46	6.46	95.6	ng	5177680	1	6.68	1082990	20.0
2-Pyrrolidinone	7.13	2.4	ng	129273	1	6.68	1082990	20.0
Hexadecane	9.66	2.4	ng	140292	3	9.72	1177380	20.0
Eicosane	10.37	2.2	ng	131703	3	9.72	1177380	20.0
Tridecanoic acid	11.75	3.5	ng	166589	4	11.20	964866	20.0
Tetradecanoic acid	12.52	2.1	ng	83580	5	13.83	809505	20.0
1-Docosene	13.69	19.3	ng	779145	5	13.83	809505	20.0
Oleic Acid	14.27	4.4	ng	178451	5	13.83	809505	20.0
1-Hexadecanesulfo...	15.70	4.0	ng	143244	6	15.20	709423	20.0
Hexacosane	16.04	2.1	ng	73370	6	15.20	709423	20.0
unknown16.58	16.58	3.9	ng	139778	6	15.20	709423	20.0
unknown17.31	17.31	2.8	ng	100885	6	15.20	709423	20.0