

Data Path : Z:\HPCHEM1\BNA F\DATA\BF050316\
 Data File : BF086651.D
 Acq On : 3 May 2016 18:22
 Operator : UM/SJ
 Sample : H2685-09
 Misc :
 ALS Vial : 41 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 26WARD-2B-CS-8(10-20FTBGS)

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-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	3.367	109	112	119	rBV	28979	64050	1.03%	0.129%
2	4.247	187	189	203	rBV	29886	87713	1.41%	0.177%
3	4.944	248	250	256	rBV	141257	250428	4.03%	0.505%
4	5.356	283	286	288	rBV	5165648	6130986	98.64%	12.354%
5	6.396	374	377	380	rBV	4334852	6215723	100.00%	12.525%
6	6.522	385	388	391	rVV	5550069	5600626	90.10%	11.286%
7	6.750	406	408	419	rBV	1383659	1312867	21.12%	2.646%
8	6.910	419	422	424	rBV	4587882	4510365	72.56%	9.089%
9	7.322	455	458	460	rBV	3525340	3815672	61.39%	7.689%
10	8.042	518	521	523	rBV	1340530	1572078	25.29%	3.168%
11	8.579	565	568	570	rVB	284833	321090	5.17%	0.647%
12	8.899	593	596	603	rBV2	108035	198536	3.19%	0.400%
13	9.116	612	615	618	rBV	4592912	4799851	77.22%	9.672%
14	9.516	646	650	652	rBV	875954	951292	15.30%	1.917%
15	9.733	665	669	671	rBV	93149	83967	1.35%	0.169%
16	9.790	671	674	676	rBV	1854995	1651151	26.56%	3.327%
17	10.225	705	712	714	rBV2	112497	198029	3.19%	0.399%
18	10.453	728	732	733	rBV	62410	101042	1.63%	0.204%
19	10.579	740	743	745	rBV	2845042	2988915	48.09%	6.023%
20	10.705	752	754	761	rVB	218286	278861	4.49%	0.562%
21	11.151	791	793	794	rBV	118034	97767	1.57%	0.197%
22	11.265	801	803	807	rBV	1088289	1285942	20.69%	2.591%
23	11.814	848	851	855	rBV2	189910	233330	3.75%	0.470%
24	12.476	907	909	912	rBV	80521	81270	1.31%	0.164%
25	12.682	925	927	931	rBV3	239574	340449	5.48%	0.686%
26	12.854	939	942	945	rBV	2880247	3030241	48.75%	6.106%
27	13.391	987	989	991	rBV	219430	191680	3.08%	0.386%
28	13.768	1019	1022	1028	rBV	165108	320726	5.16%	0.646%
29	13.905	1031	1034	1036	rBV	726909	1078748	17.36%	2.174%
30	14.008	1041	1043	1047	rVV3	50444	99413	1.60%	0.200%
31	14.077	1047	1049	1053	rVB	37562	63630	1.02%	0.128%
32	14.385	1074	1076	1078	rBV	57635	72661	1.17%	0.146%
33	14.660	1097	1100	1106	rBV	448509	552266	8.88%	1.113%
34	14.911	1119	1122	1125	rBV	33155	70228	1.13%	0.142%

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ALS Vial : 41 Sample Multiplier: 1

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BNA_F
ClientSampleId :
26WARD-2B-CS-8(10-20FTBGS)

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

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35	15.300	1153	1156	1158	rBV	608812	906438	14.58%	1.827%
36	16.191	1231	1234	1236	rBV	29573	68040	1.09%	0.137%

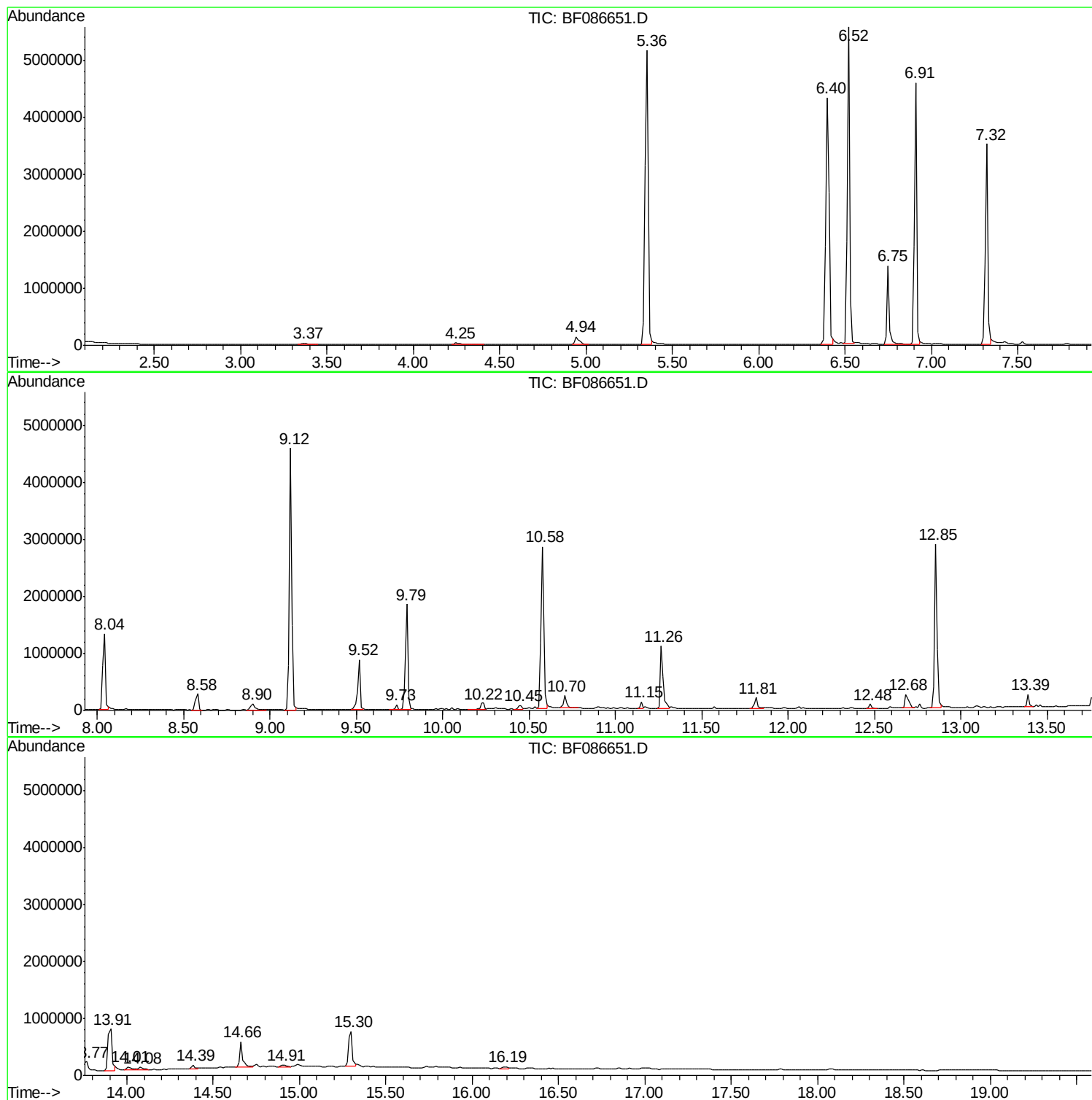
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF042716.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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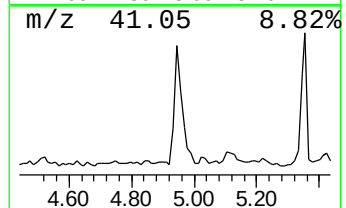
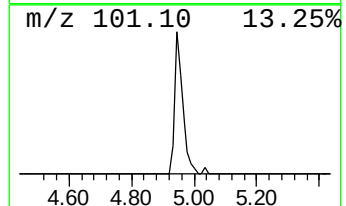
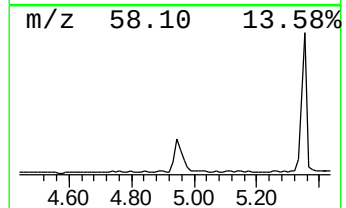
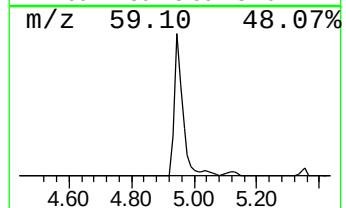
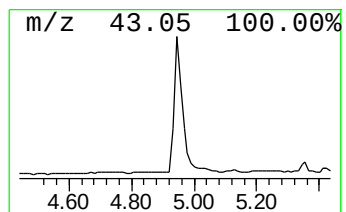
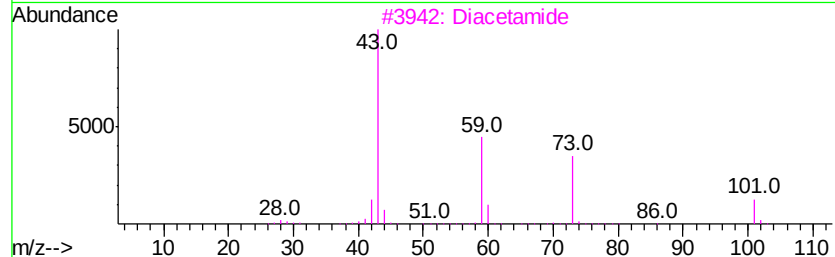
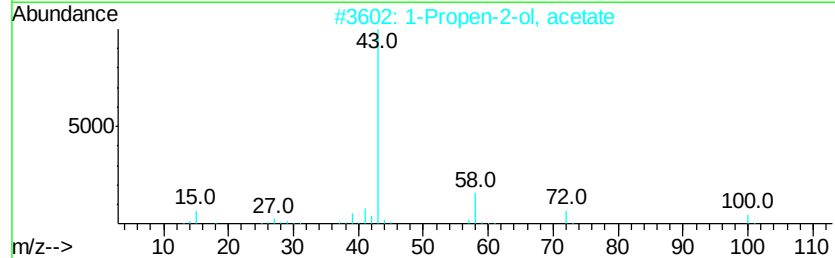
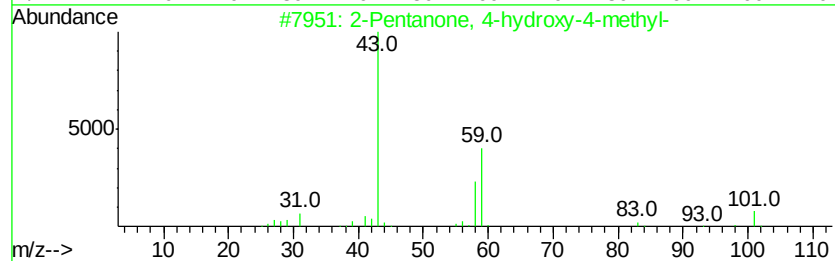
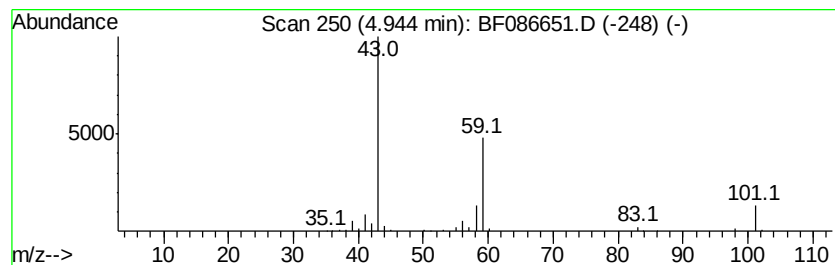
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.94	3.81 ng	250428	1,4-Dichlorobenzene-d4	6.75

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2		1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10
3		Diacetamide	101	C4H7NO2	000625-77-4	9
4		Butane	58	C4H10	000106-97-8	9
5		Acetic acid, 2-propenyl ester	100	C5H8O2	000591-87-7	9



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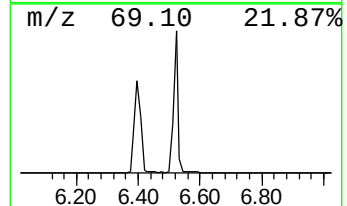
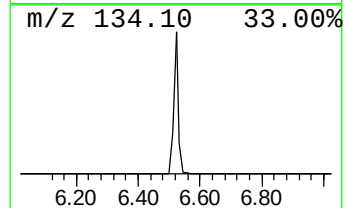
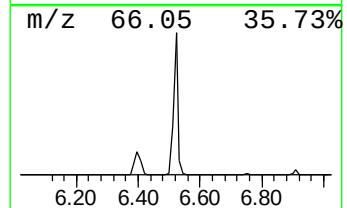
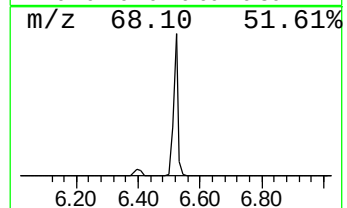
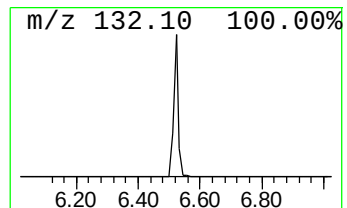
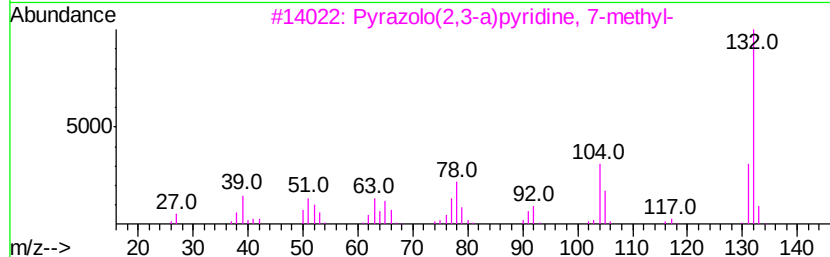
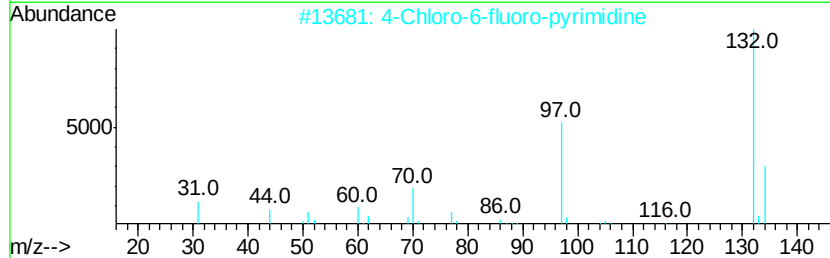
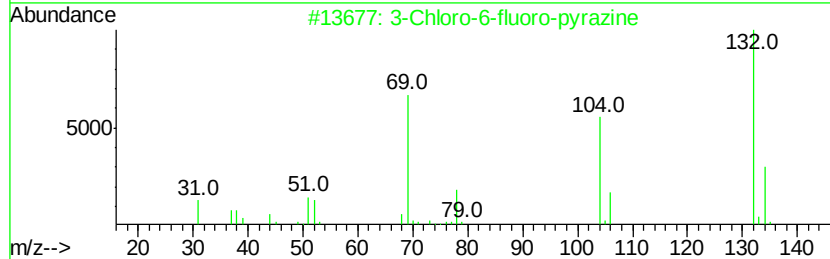
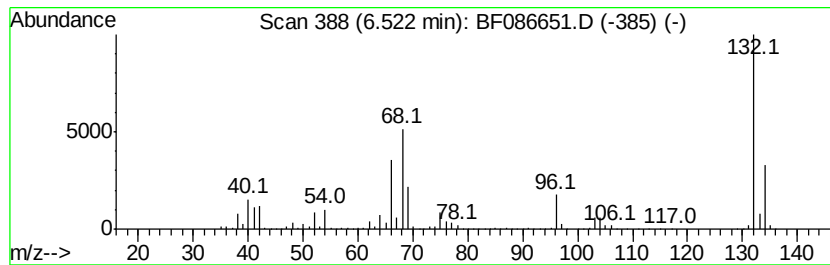
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown6.52 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.52	85.32 ng	5600630	1,4-Dichlorobenzene-d4	6.75

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	27
3		Pyrazolo(2,3-a)pyridine, 7-methyl-	132	C8H8N2	034760-58-2	22
4		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
5		Tranlycypromine-propionyl	189	C12H15NO	1000123-86-3	22



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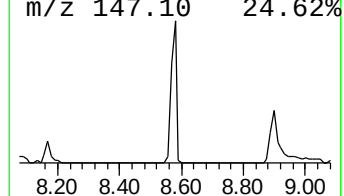
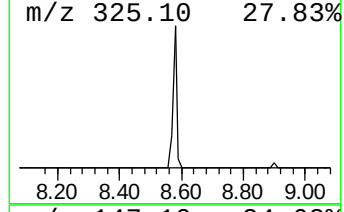
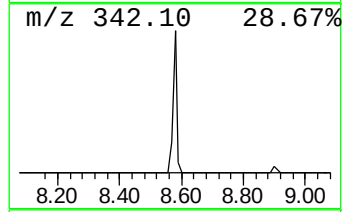
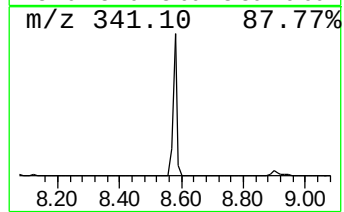
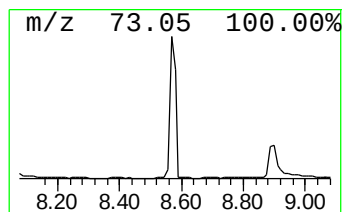
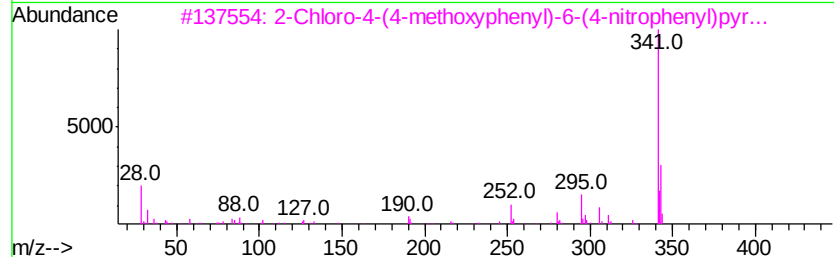
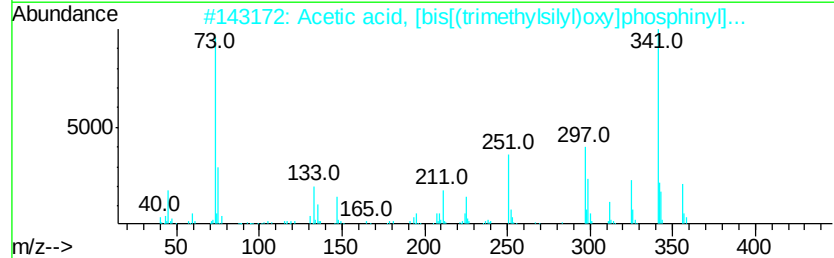
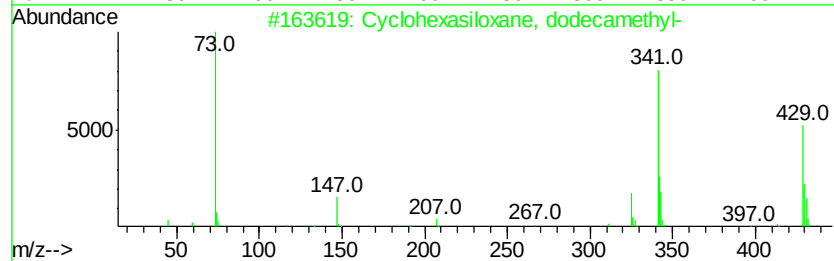
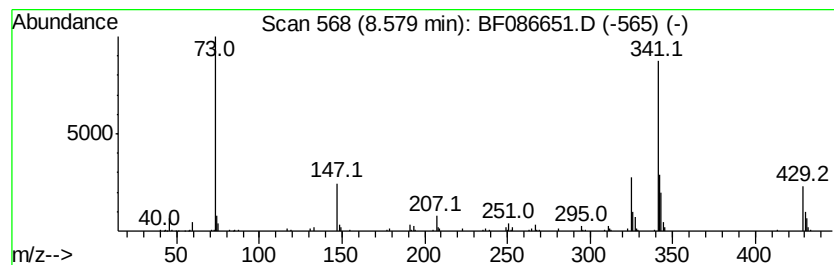
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Cyclohexasiloxane, dodecane... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.58	4.08 ng	321090	Naphthalene-d8	8.04

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexasiloxane, dodecamethyl-	444	C12H36O6Si6	000540-97-6	91
2		Acetic acid, [bis[(trimethylsilyl)oxy]phosphiny]...	356	C11H29O5PSi3	053044-27-2	53
3		2-Chloro-4-(4-methoxyphenyl)-6-(4-nitrophenyl)pyr...	341	C17H12ClN3O3	063673-76-7	38
4		Benzothiophene, 5-chloro-3-methyl-	341	C18H12ClNS2	1000270-14-7	17
5		(3-Amino-6-thiophen-2-yl-thieno[3,2-b]pyridine)	342	C16H10N2OS3	1000296-68-0	13



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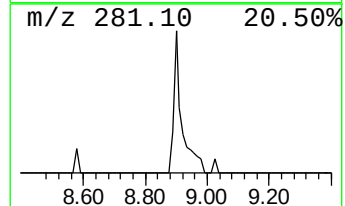
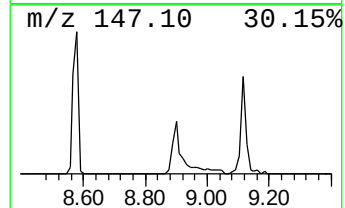
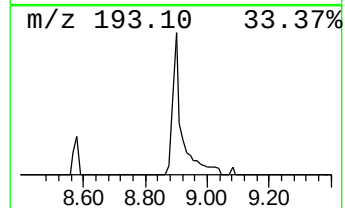
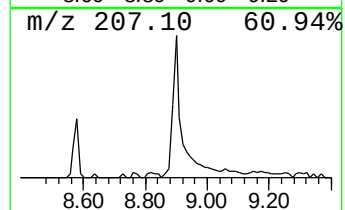
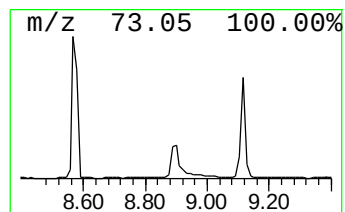
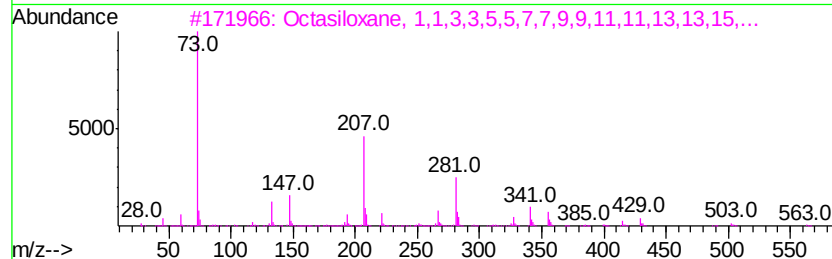
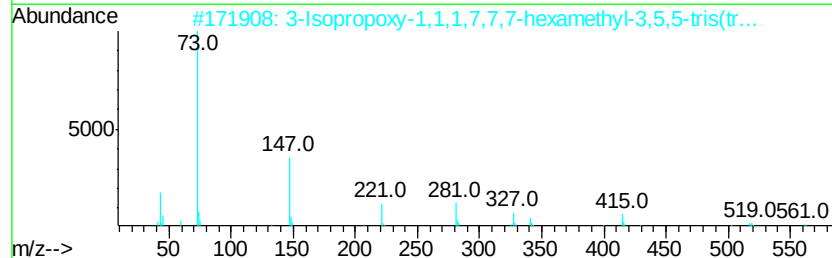
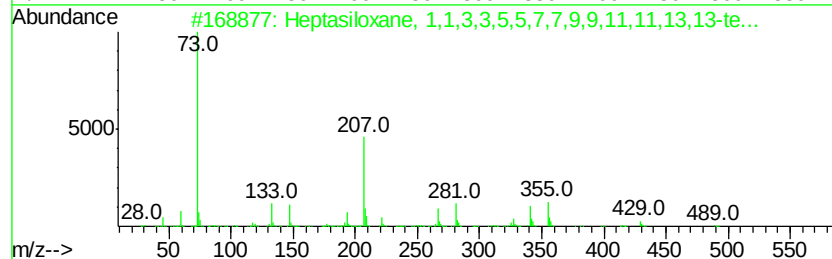
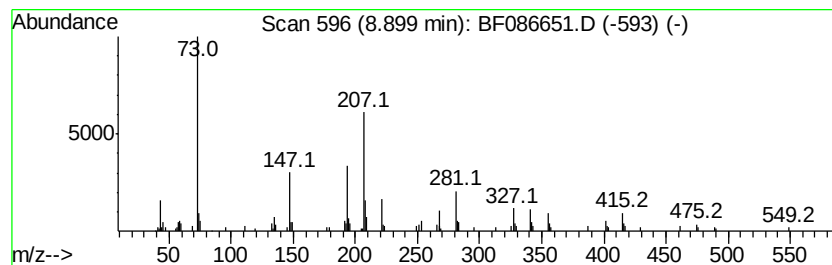
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 5 unknown8.90 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.90	2.53 ng	198536	Napthalene-d8	8.04

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptasiloxane, 1,1,3,3,5,5,7,7,9...	504	C14H44O6Si7	019095-23-9	47
2		3-Isopropoxy-1,1,1,7,7,7-hexamet...	576	C18H52O7Si7	071579-69-6	41
3		Octasiloxane, 1,1,3,3,5,5,7,7,9...	578	C16H50O7Si8	019095-24-0	37
4		Benzo[h]quinoline, 2,4-dimethyl-	207	C15H13N	000605-67-4	25
5		5,5'-Di(ethoxycarbonyl)-3,3'-dim...	402	C23H34N2O4	102586-97-0	25



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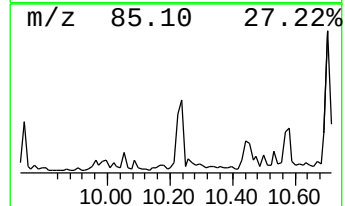
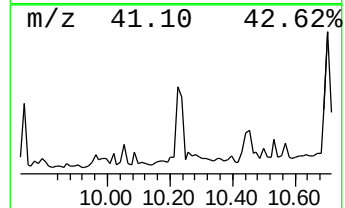
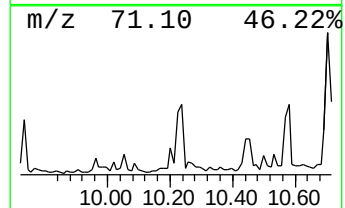
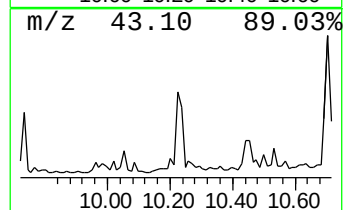
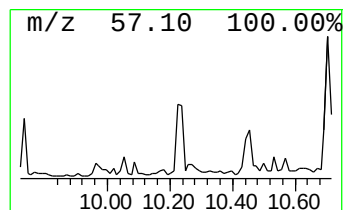
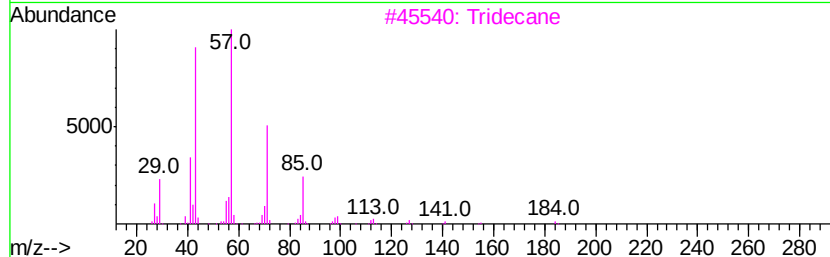
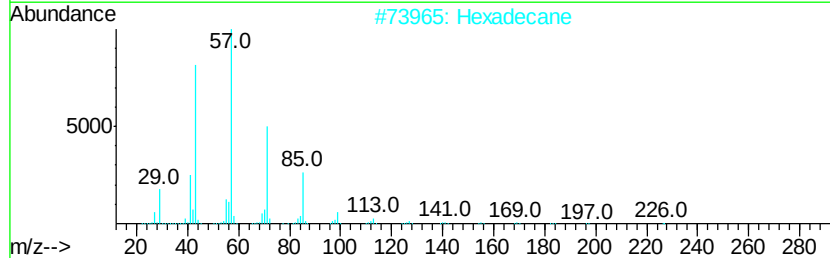
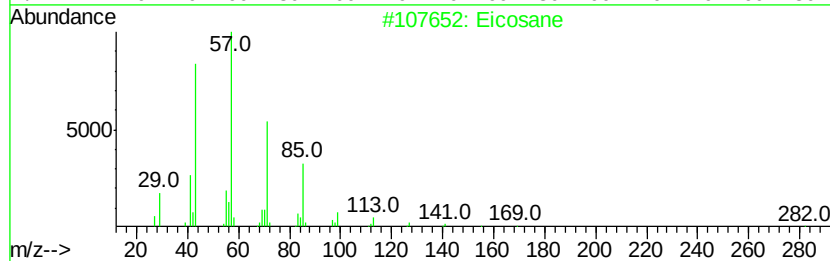
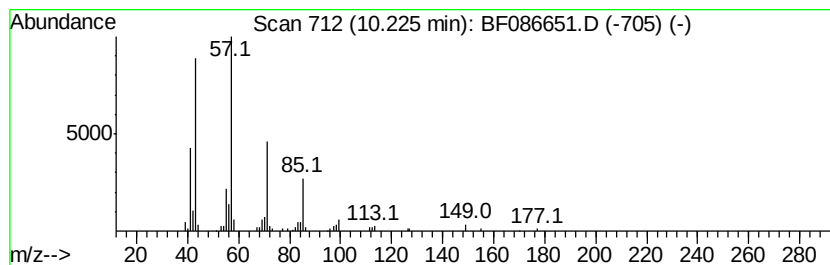
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Peak Number 6 Eicosane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.22	2.40 ng	198029	Acenaphthene-d10	9.79
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Eicosane		282 C20H42	000112-95-8 91
2	Hexadecane		226 C16H34	000544-76-3 91
3	Tridecane		184 C13H28	000629-50-5 90
4	Octadecane		254 C18H38	000593-45-3 86
5	Pentadecane		212 C15H32	000629-62-9 86



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Acq On : 3 May 2016 18:22
Operator : UM/SJ
Sample : H2685-09
Misc :
ALS Vial : 41 Sample Multiplier: 1

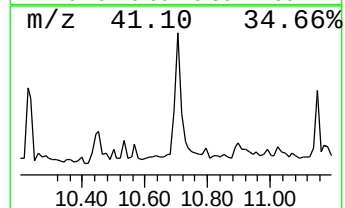
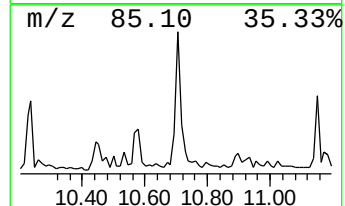
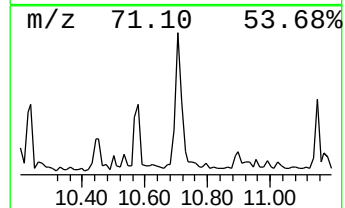
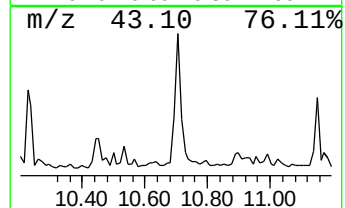
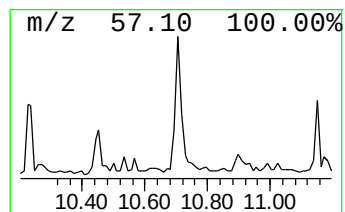
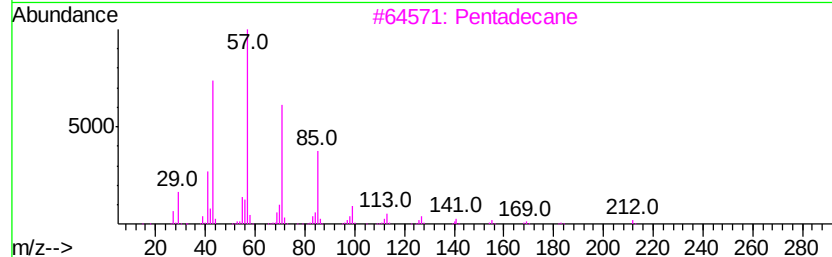
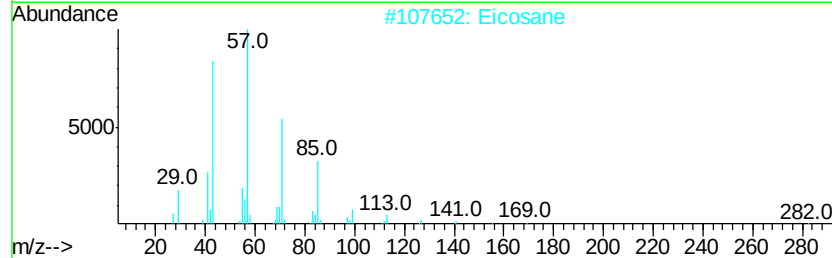
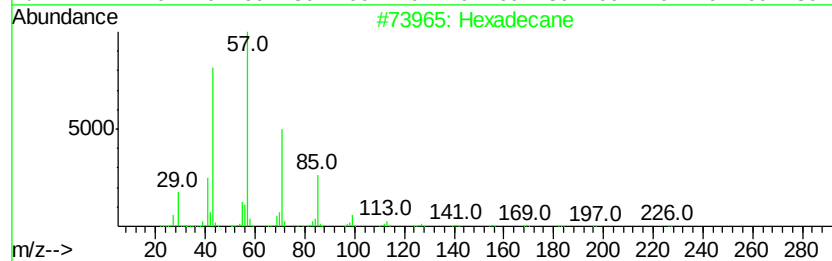
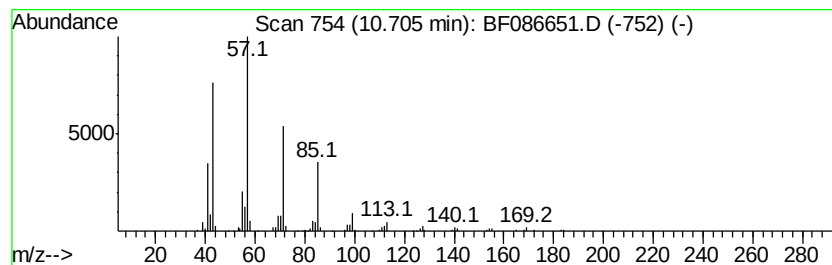
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ClientSampleId :
26WARD-2B-CS-8(10-20FTBGS)

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 7 Hexadecane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.70	4.34 ng	278861	Phenanthrene-d10	11.26	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane	226	C16H34	000544-76-3	91
2	Eicosane	282	C20H42	000112-95-8	91
3	Pentadecane	212	C15H32	000629-62-9	90
4	Dodecane, 2-methyl-6-propyl-	226	C16H34	055045-08-4	90
5	Tridecane	184	C13H28	000629-50-5	90



Data Path : Z:\HPCHEM1\BNA F\DATA\BF050316\
Data File : BF086651.D
Acq On : 3 May 2016 18:22
Operator : UM/SJ
Sample : H2685-09
Misc :
ALS Vial : 41 Sample Multiplier: 1

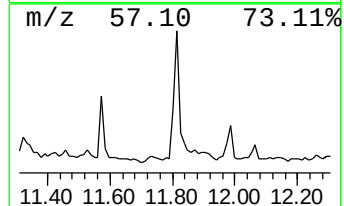
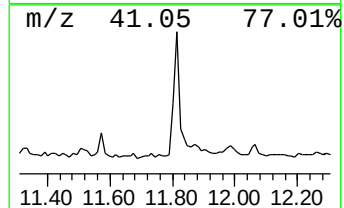
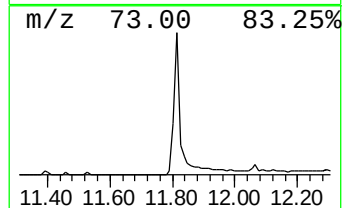
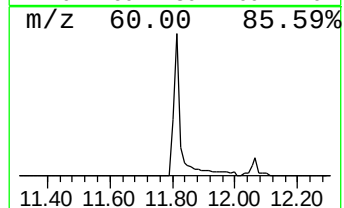
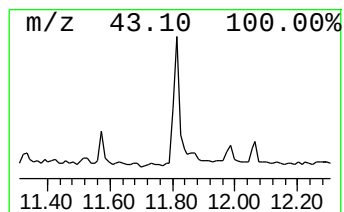
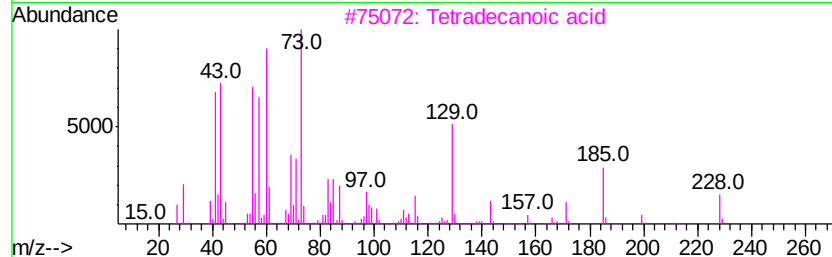
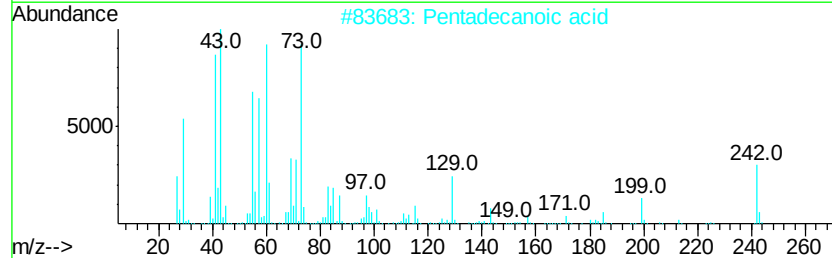
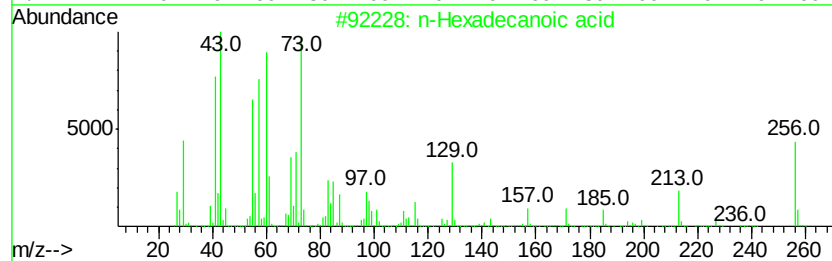
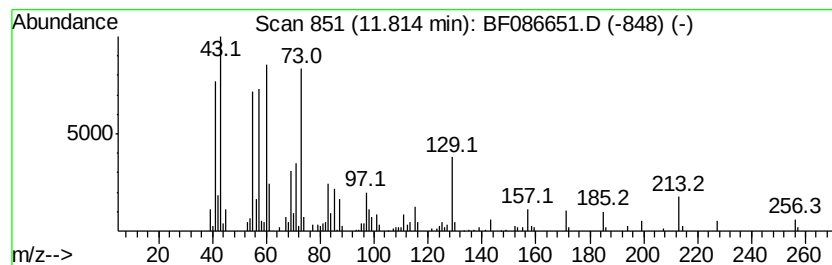
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ClientSampleId :
26WARD-2B-CS-8(10-20FTBGS)

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 n-Hexadecanoic acid Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD		R.T.
11.81	3.63 ng	233330	Phenanthrene-d10		11.26
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	94
2	Pentadecanoic acid	242	C15H30O2	001002-84-2	80
3	Tetradecanoic acid	228	C14H28O2	000544-63-8	72
4	Tridecanoic acid	214	C13H26O2	000638-53-9	64
5	n-Decanoic acid	172	C10H20O2	000334-48-5	53



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF050316\
Data File : BF086651.D
Acq On : 3 May 2016 18:22
Operator : UM/SJ
Sample : H2685-09
Misc :
ALS Vial : 41 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
26WARD-2B-CS-8(10-20FTBGS)

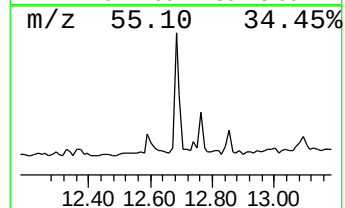
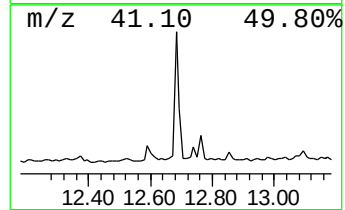
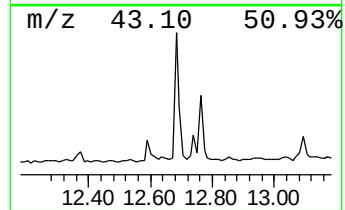
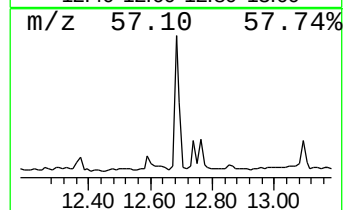
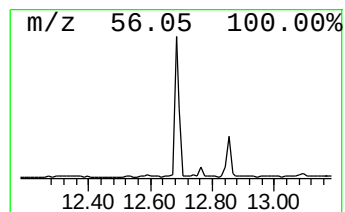
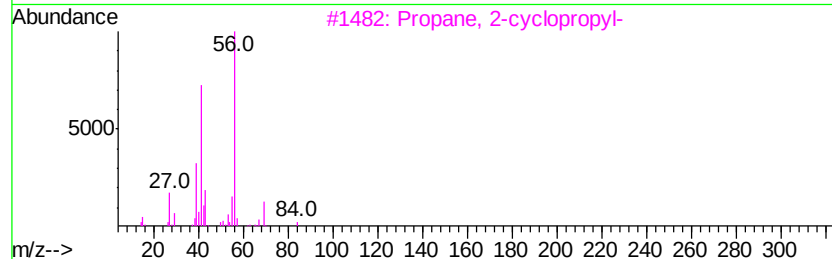
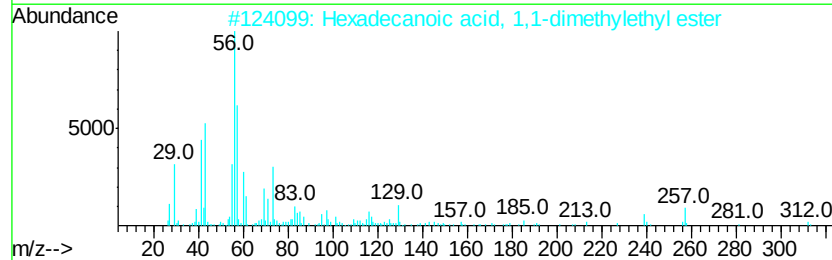
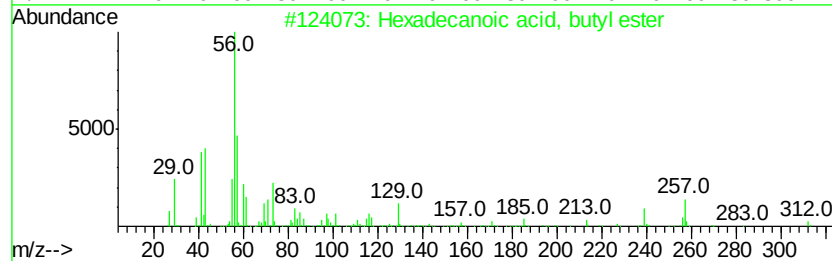
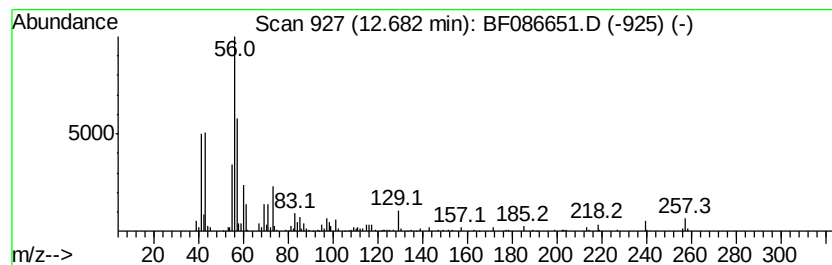
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Hexadecanoic acid, butyl ester Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.68	6.31 ng	340449	Chrysene-d12	13.91

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecanoic acid, butyl ester	312	C20H40O2	000111-06-8	91
2		Hexadecanoic acid, 1,1-dimethyle...	312	C20H40O2	031158-91-5	68
3		Propane, 2-cyclopropyl-	84	C6H12	003638-35-5	43
4		1-Hexene, 2-methyl-	98	C7H14	006094-02-6	38
5		Hexadecanoic acid, 2-methylpropy...	312	C20H40O2	000110-34-9	38



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF050316\
Data File : BF086651.D
Acq On : 3 May 2016 18:22
Operator : UM/SJ
Sample : H2685-09
Misc :
ALS Vial : 41 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
26WARD-2B-CS-8(10-20FTBGS)

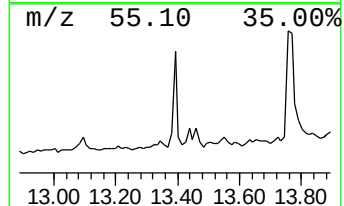
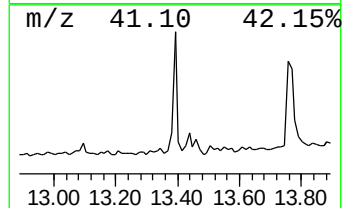
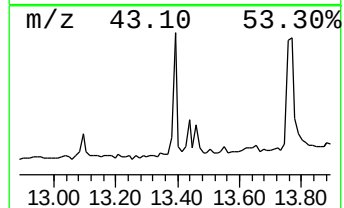
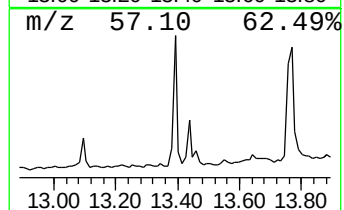
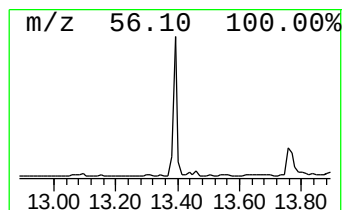
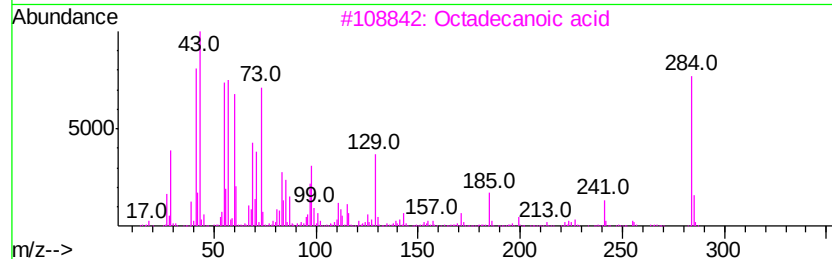
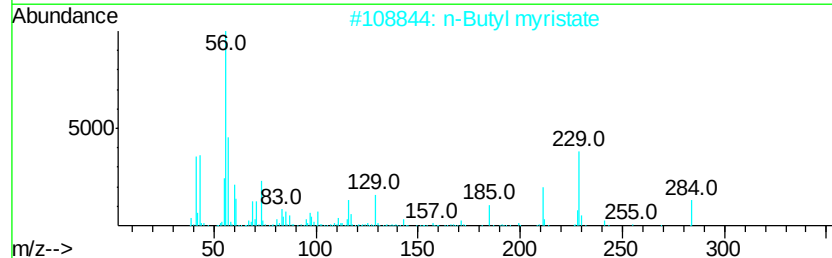
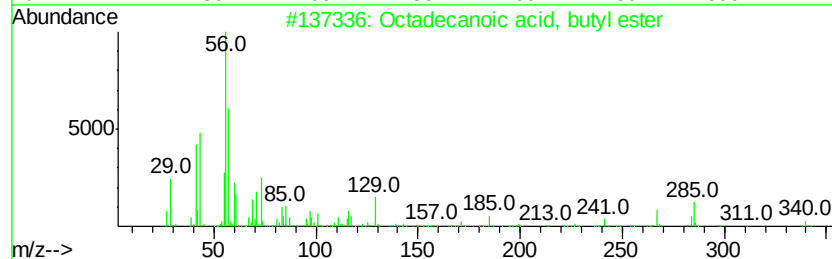
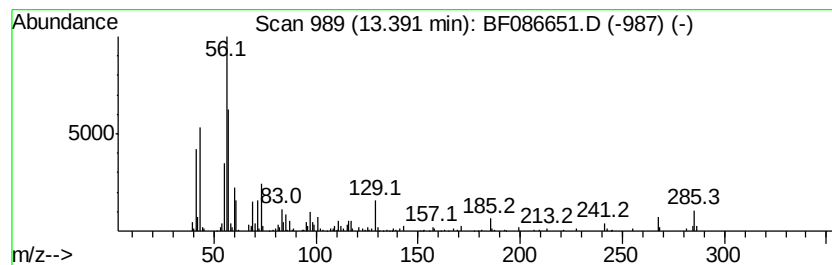
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Octadecanoic acid, butyl ester Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.39	3.55 ng	191680	Chrysene-d12	13.91

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid, butyl ester	340	C22H44O2	000123-95-5	91
2			n-Butyl myristate	284	C18H36O2	000110-36-1	78
3			Octadecanoic acid	284	C18H36O2	000057-11-4	38
4			Urea, N,N'-di-2-propenyl-	140	C7H12N2O	001801-72-5	38
5			Cyclopentanone, 2,2,5,5-tetramet...	140	C9H16O	004541-35-9	38



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF050316\
Data File : BF086651.D
Acq On : 3 May 2016 18:22
Operator : UM/SJ
Sample : H2685-09
Misc :
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ClientSampleId :
26WARD-2B-CS-8(10-20FTBGS)

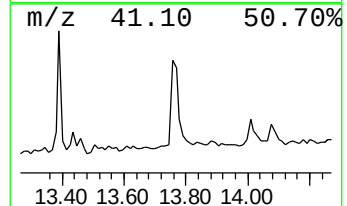
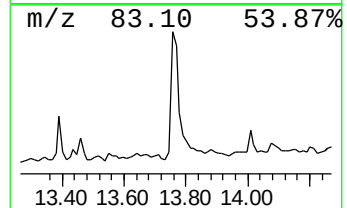
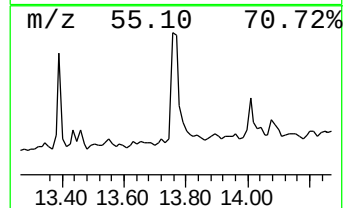
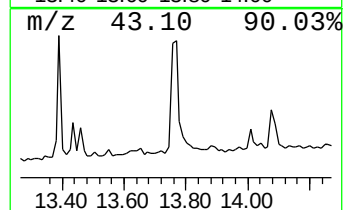
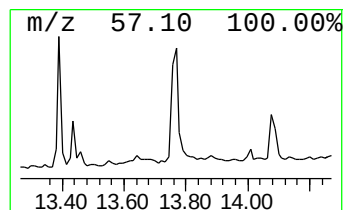
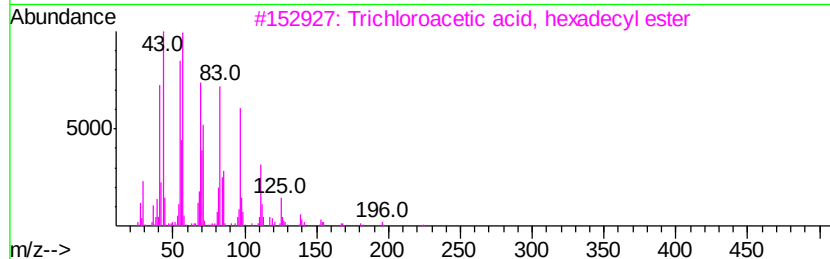
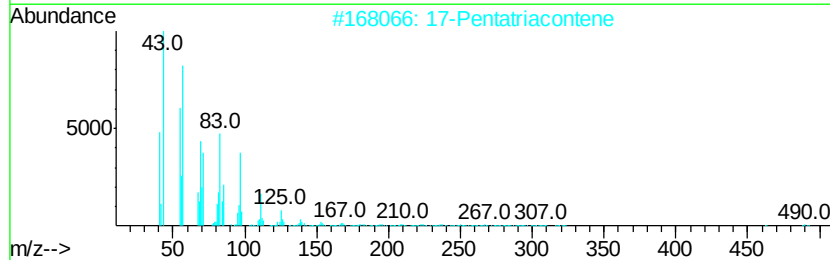
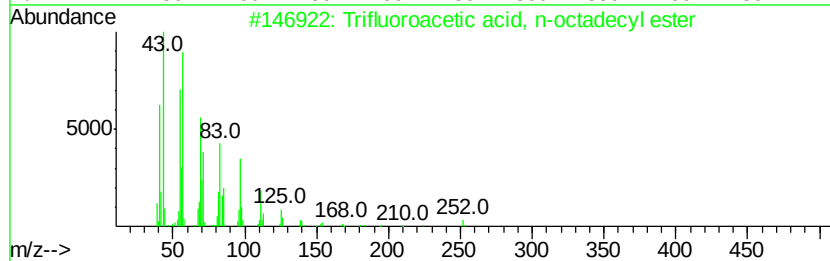
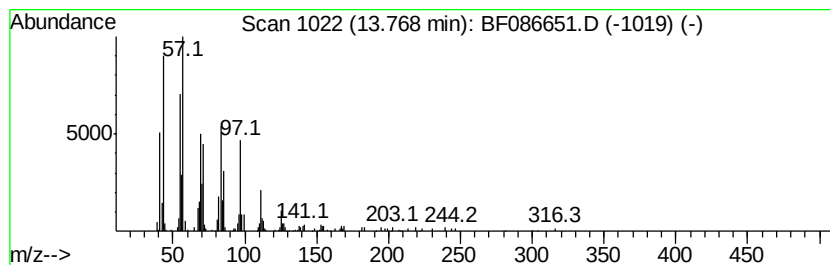
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Trifluoroacetic acid, n-oct... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.77	5.95 ng	320726	Chrysene-d12	13.91

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Trifluoroacetic acid, n-octadecyl...	366	C20H37F3O2	079392-43-1	91
2		17-Pentatriacontene	491	C35H70	006971-40-0	91
3		Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	90
4		Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	87
5		Pentafluoropropionic acid, hepta...	402	C20H35F5O2	1000283-04-2	87



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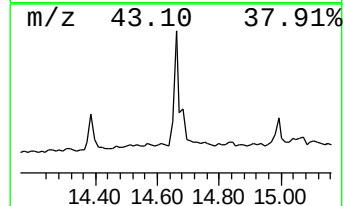
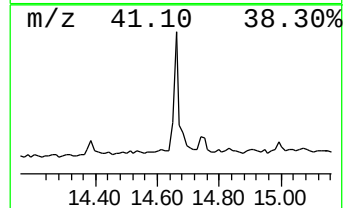
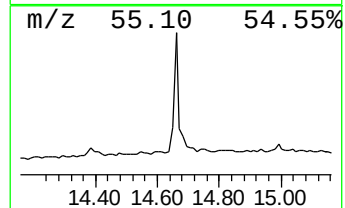
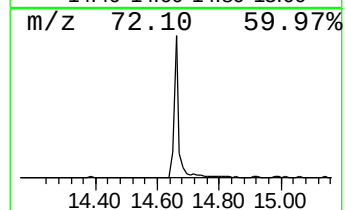
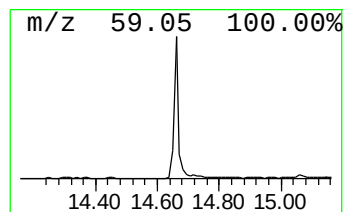
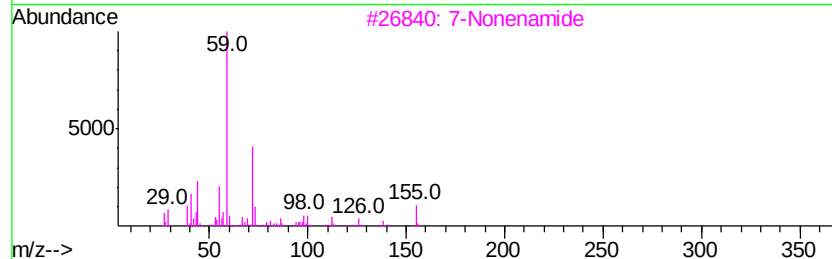
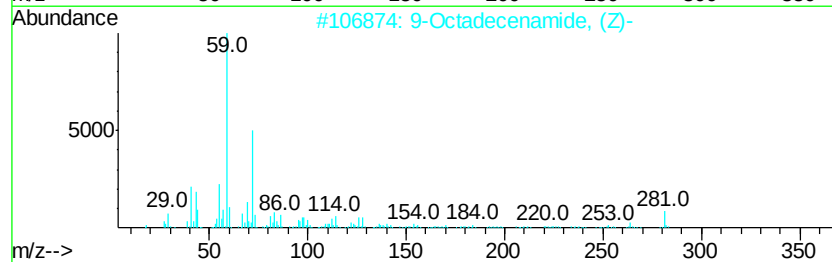
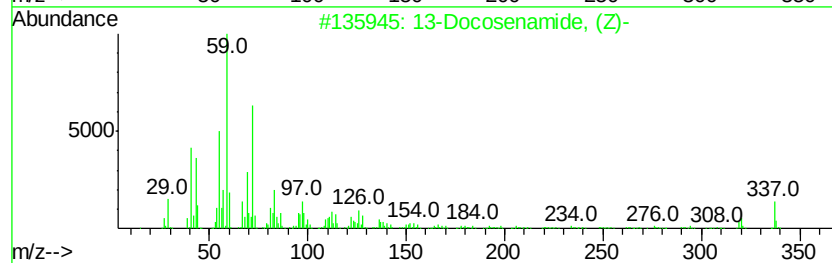
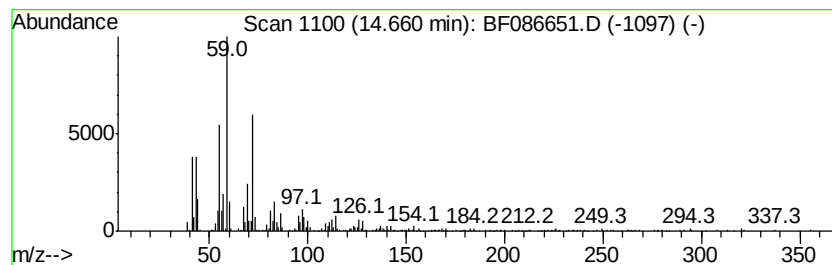
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 13-Docosenamide, (Z)- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.66	12.19 ng	552266	Perylene-d12	15.30

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	13-Docosenamide, (Z)-	337	C22H43NO	000112-84-5	95
2		9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	87
3		7-Nonenamide	155	C9H17NO	090949-53-4	64
4		Octanamide	143	C8H17NO	000629-01-6	59
5		Hexadecanamide	255	C16H33NO	000629-54-9	59



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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
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unknown6.52	6.52	85.3	ng	5600630	1	6.75	1312870	20.0
Cyclohexasiloxane...	8.58	4.1	ng	321090	2	8.04	1572080	20.0
unknown8.90	8.90	2.5	ng	198536	2	8.04	1572080	20.0
Eicosane	10.22	2.4	ng	198029	3	9.79	1651150	20.0
Hexadecane	10.70	4.3	ng	278861	4	11.26	1285940	20.0
n-Hexadecanoic acid	11.81	3.6	ng	233330	4	11.26	1285940	20.0
Hexadecanoic acid...	12.68	6.3	ng	340449	5	13.91	1078750	20.0
Octadecanoic acid...	13.39	3.5	ng	191680	5	13.91	1078750	20.0
Trifluoroacetic a...	13.77	6.0	ng	320726	5	13.91	1078750	20.0
13-Docosenamide, ...	14.66	12.2	ng	552266	6	15.30	906438	20.0