

Data Path : Z:\HPCHEM1\BNA F\DATA\BF050316\
 Data File : BF086648.D
 Acq On : 3 May 2016 16:57
 Operator : UM/SJ
 Sample : H2685-06
 Misc :
 ALS Vial : 38 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 26WARD-2B-CS-6(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	3.356	107	111	118	rBV	40524	89119	1.71%	0.180%
2	4.236	186	188	193	rBV	39552	75868	1.46%	0.153%
3	4.933	247	249	256	rBV	142935	231660	4.46%	0.467%
4	5.344	282	285	288	rBV	3626796	4997283	96.16%	10.082%
5	6.396	374	377	380	rBV	4374931	4962513	95.49%	10.012%
6	6.522	385	388	390	rBV	5218541	5197052	100.00%	10.485%
7	6.750	406	408	419	rBV	1414090	1320751	25.41%	2.665%
8	6.910	419	422	424	rBV	4125045	4073716	78.39%	8.218%
9	7.322	455	458	460	rBV	3339852	3488914	67.13%	7.039%
10	7.425	466	467	474	rVB	48954	68321	1.31%	0.138%
11	7.527	474	476	479	rVB	60518	62100	1.19%	0.125%
12	8.042	518	521	523	rBV	1379941	1604223	30.87%	3.236%
13	8.579	565	568	570	rBV	265784	286697	5.52%	0.578%
14	8.899	593	596	602	rBV2	128252	221086	4.25%	0.446%
15	9.116	612	615	618	rBV	4440314	4387495	84.42%	8.852%
16	9.516	647	650	652	rBV	841177	786235	15.13%	1.586%
17	9.733	667	669	671	rBV	122996	102509	1.97%	0.207%
18	9.790	671	674	676	rBV	1901105	1712105	32.94%	3.454%
19	10.236	709	713	714	rBV	160995	246147	4.74%	0.497%
20	10.453	729	732	735	rBV	86288	151477	2.91%	0.306%
21	10.579	740	743	745	rBV	2870485	2922547	56.23%	5.896%
22	10.705	752	754	760	rVB	307926	374663	7.21%	0.756%
23	10.899	768	771	773	rBV2	38363	67499	1.30%	0.136%
24	11.151	791	793	794	rBV	158159	126944	2.44%	0.256%
25	11.173	794	795	799	rVB	38748	56562	1.09%	0.114%
26	11.265	801	803	807	rBV	985118	1571395	30.24%	3.170%
27	11.345	807	810	812	rVB2	64797	92318	1.78%	0.186%
28	11.814	848	851	853	rBV2	708509	685024	13.18%	1.382%
29	11.882	855	857	862	rVB2	81121	141528	2.72%	0.286%
30	12.065	870	873	874	rBV	65641	63438	1.22%	0.128%
31	12.476	907	909	912	rBV	660499	607604	11.69%	1.226%
32	12.591	918	919	921	rBV	47866	62285	1.20%	0.126%
33	12.705	926	929	931	rBV	543923	696707	13.41%	1.406%
34	12.762	931	934	938	rVB3	48777	75947	1.46%	0.153%

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35	12.854	940	942	945	rVV	2656313	2808164	54.03%	5.665%
36	12.922	946	948	954	rVB3	46758	89446	1.72%	0.180%
37	13.037	954	958	960	rVB	48980	91683	1.76%	0.185%
38	13.094	960	963	965	rBV2	35022	61012	1.17%	0.123%
39	13.139	965	967	968	rBV	47830	63972	1.23%	0.129%
40	13.391	987	989	991	rBV	79214	82850	1.59%	0.167%
41	13.654	1009	1012	1013	rBV2	31301	52728	1.01%	0.106%
42	13.757	1019	1021	1025	rBV2	173230	307574	5.92%	0.621%
43	13.905	1030	1034	1041	rBV2	913541	1764513	33.95%	3.560%
44	14.008	1041	1043	1045	rVV	51267	55309	1.06%	0.112%
45	14.282	1065	1067	1069	rBV	34108	58306	1.12%	0.118%
46	14.385	1073	1076	1082	rVV5	41252	137508	2.65%	0.277%
47	14.660	1098	1100	1106	rBV2	104527	142678	2.75%	0.288%
48	14.751	1106	1108	1113	rVB2	95901	135590	2.61%	0.274%
49	14.900	1119	1121	1127	rBV	210976	467188	8.99%	0.943%
50	15.174	1143	1145	1148	rVB	106996	156098	3.00%	0.315%
51	15.231	1148	1150	1153	rVB	121596	175460	3.38%	0.354%
52	15.300	1153	1156	1158	rBV	682179	996100	19.17%	2.010%
53	16.614	1269	1271	1274	rBV2	57217	123006	2.37%	0.248%
54	17.014	1303	1306	1313	rVB	69090	188836	3.63%	0.381%

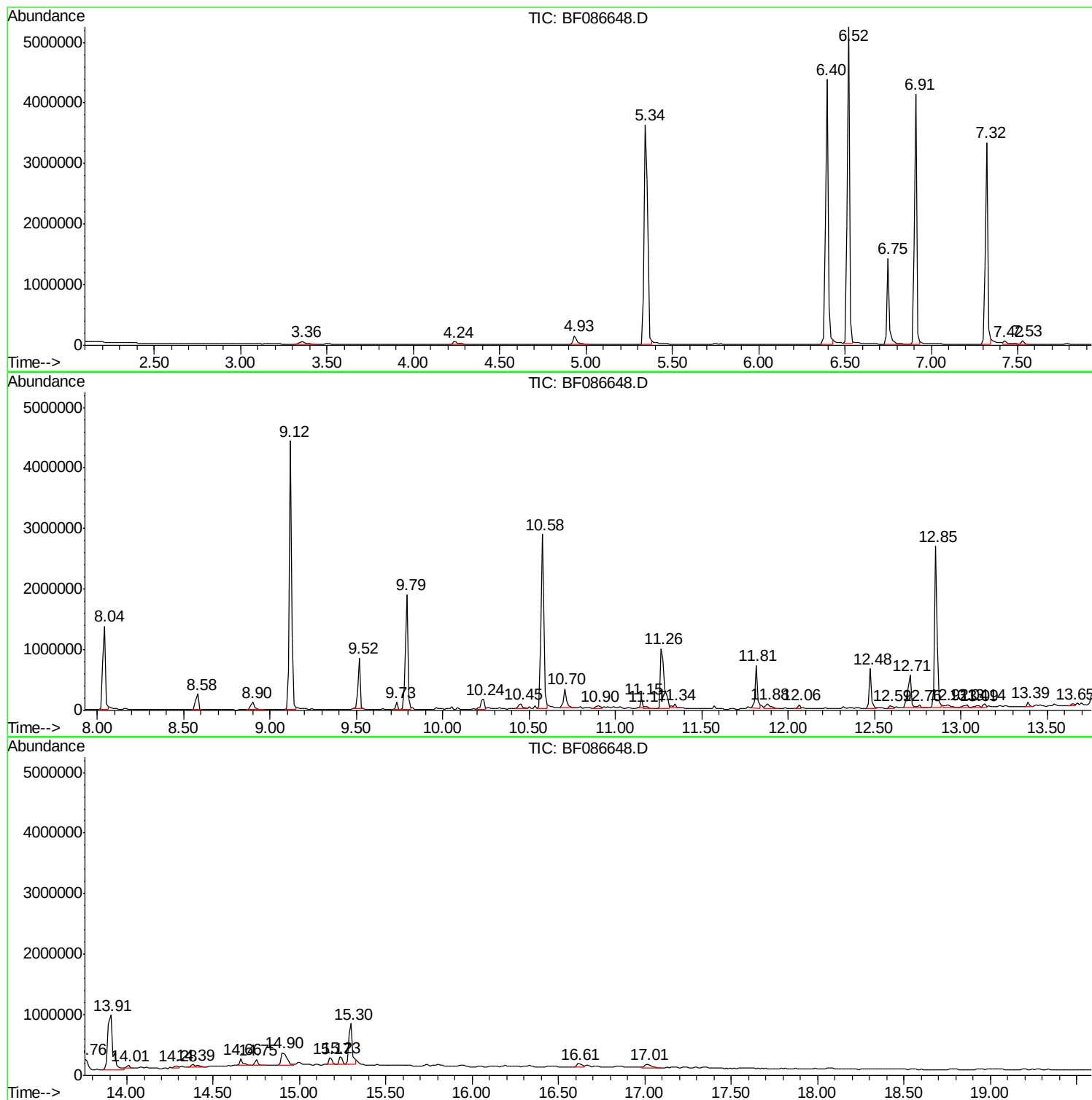
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF042716.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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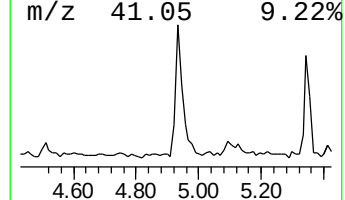
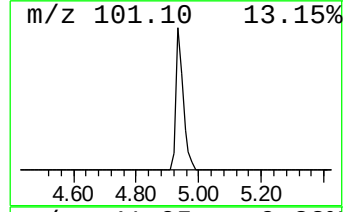
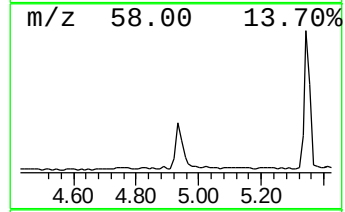
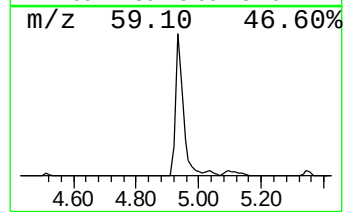
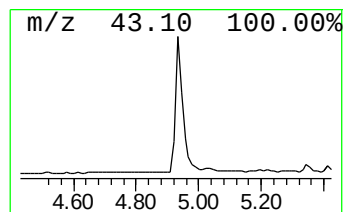
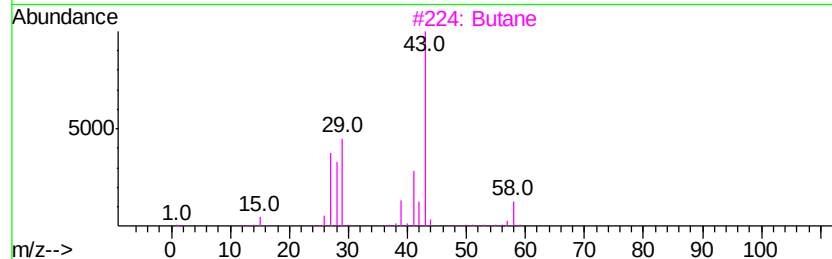
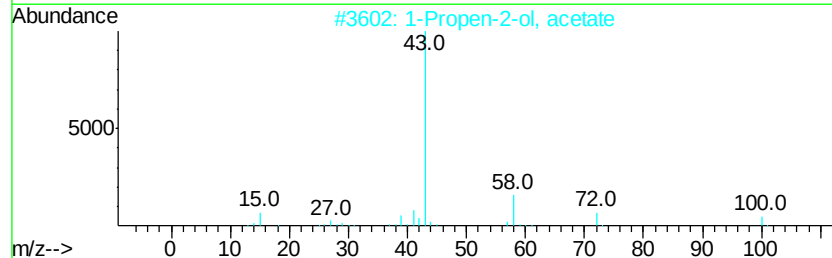
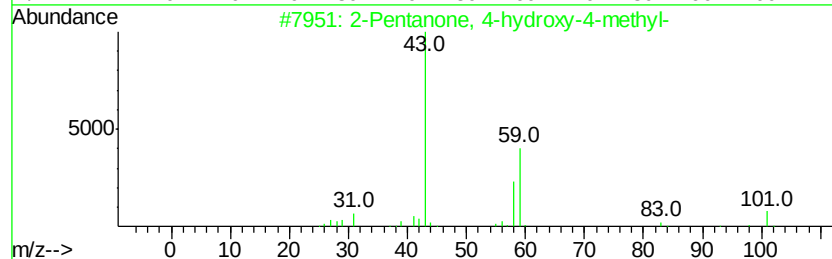
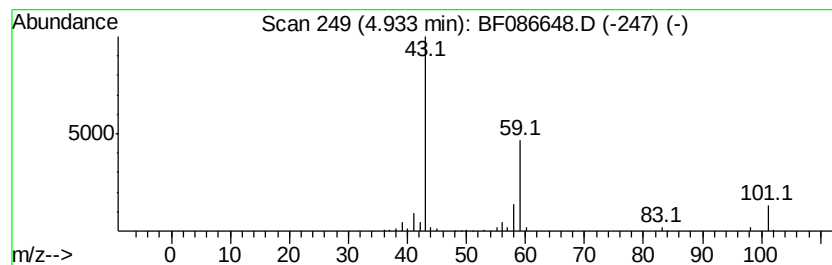
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF042716.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.93	3.51 ng	231660	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	64
2		1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10
3		Butane	58	C4H10	000106-97-8	9
4		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
5		Diacetamide	101	C4H7NO2	000625-77-4	9



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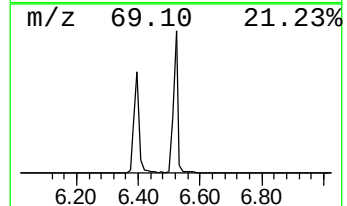
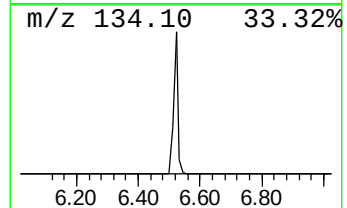
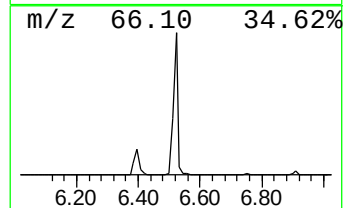
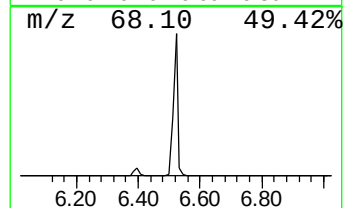
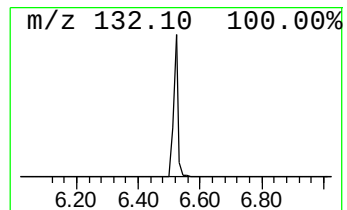
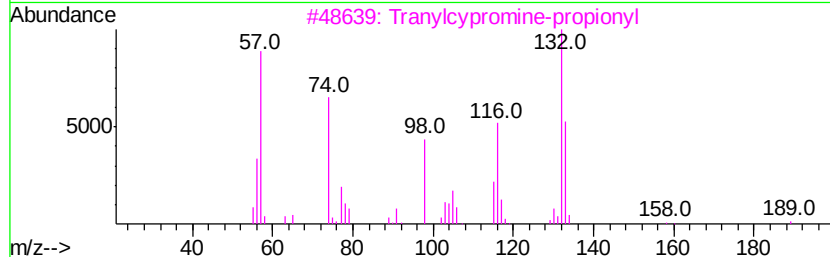
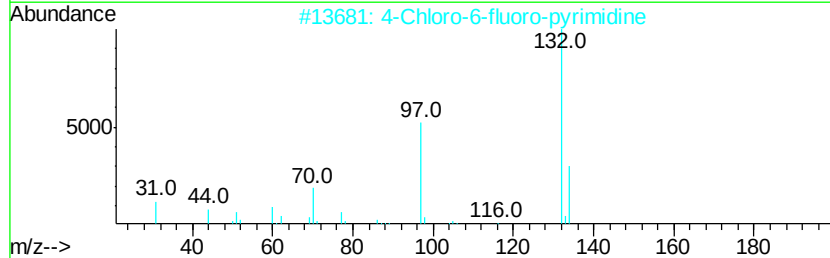
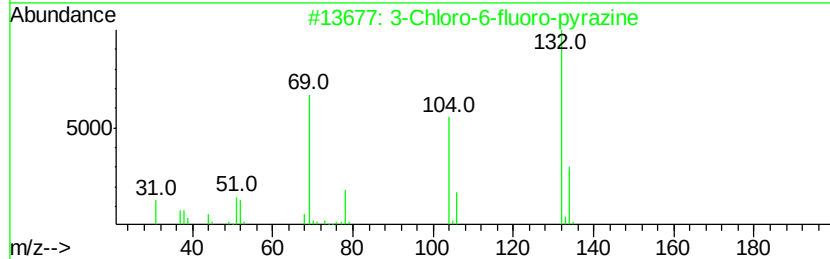
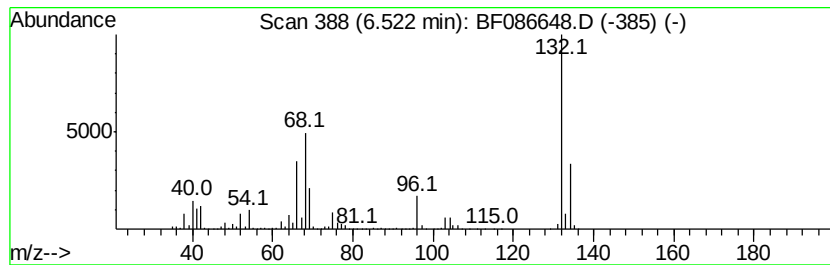
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TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown6.52 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.52	78.70 ng	5197050	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		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	22
4		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
5		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16



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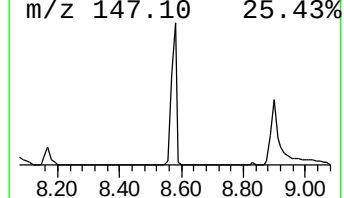
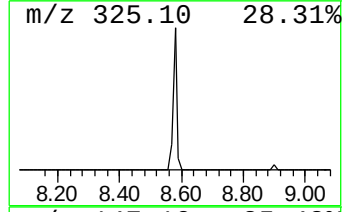
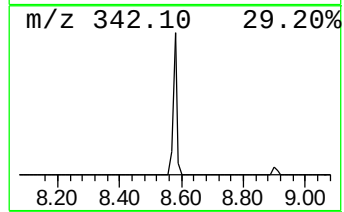
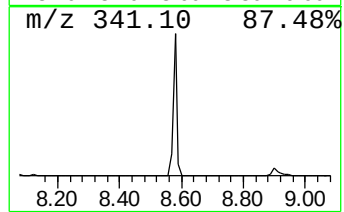
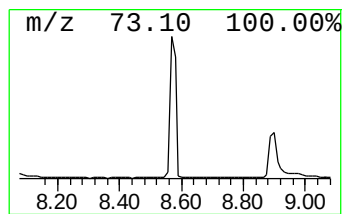
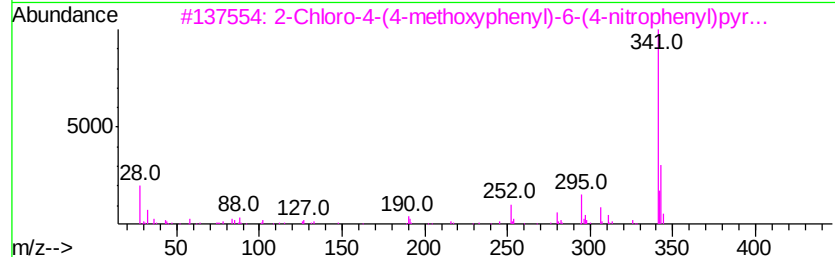
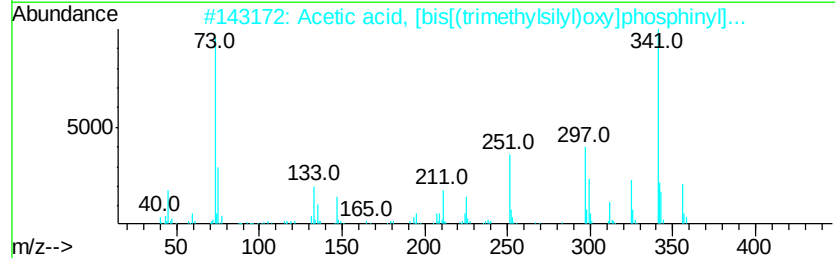
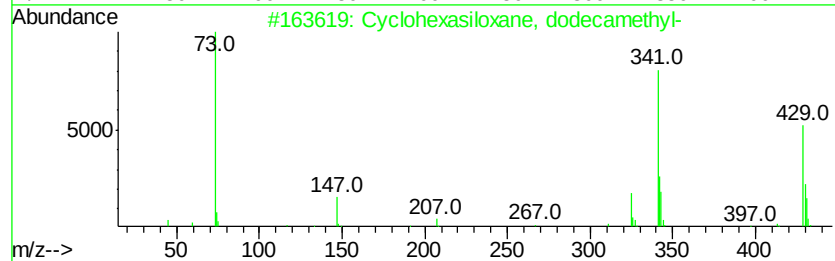
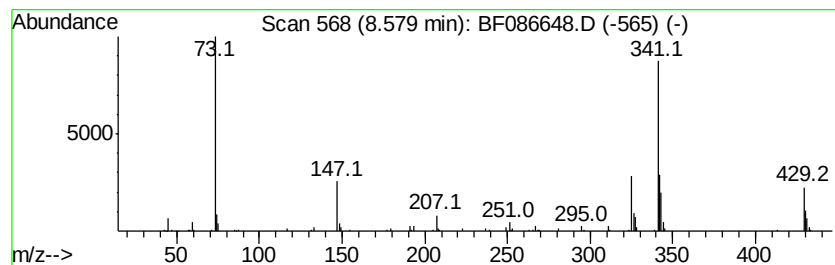
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 4 Cyclohexasiloxane, dodecane... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.58	3.57 ng	286697	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	43
4		5-Amino-7-(4-methoxy-phenyl)-2-pyr...	341	C20H15N5O	1000294-65-4	32
5		Benzothiophene, 5-chloro-3-methy...	341	C18H12ClNS2	1000270-14-7	17



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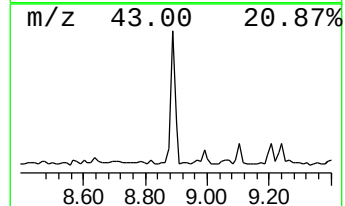
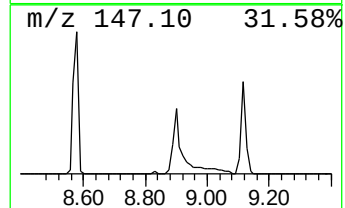
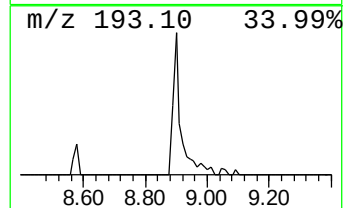
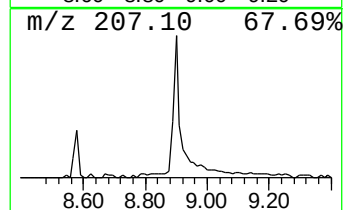
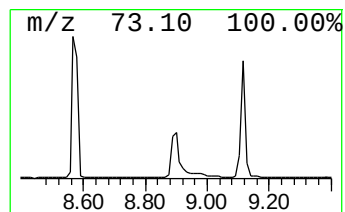
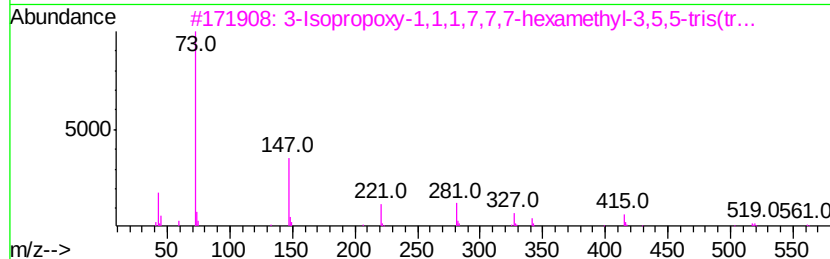
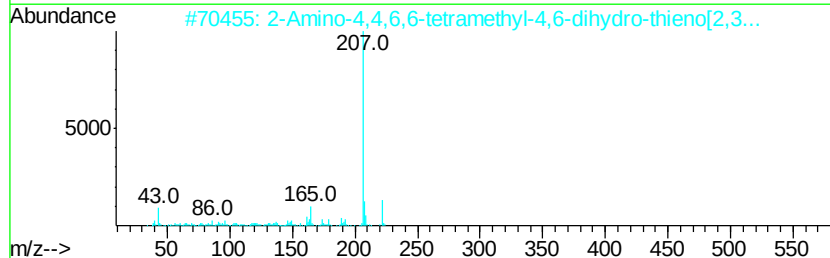
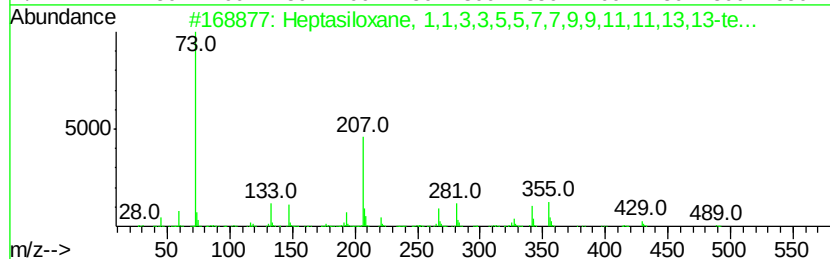
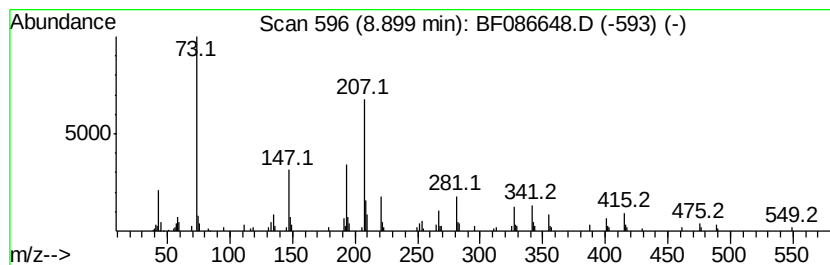
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Peak Number 5 unknown8.90 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.90	2.76 ng	221086	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	38
2		2-Amino-4,4,6,6-tetramethyl-4,6-...	222	C11H14N2OS	1000275-36-2	30
3		3-Isopropoxy-1,1,1,7,7,7-hexamet...	576	C18H52O7Si7	071579-69-6	30
4		Stannane, tetrapropyl-	292	C12H28Sn	002176-98-9	27
5		Tetrasiloxane, decamethyl-	310	C10H30O3Si4	000141-62-8	27



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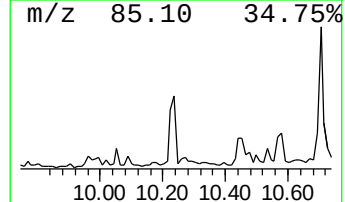
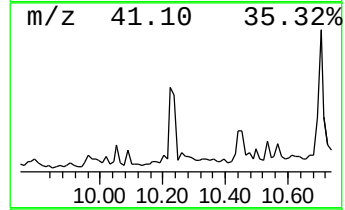
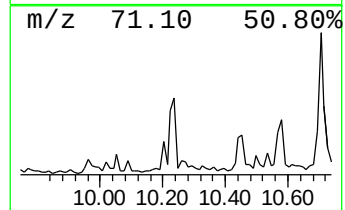
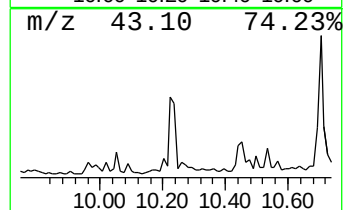
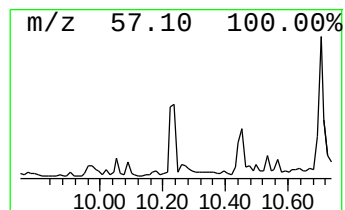
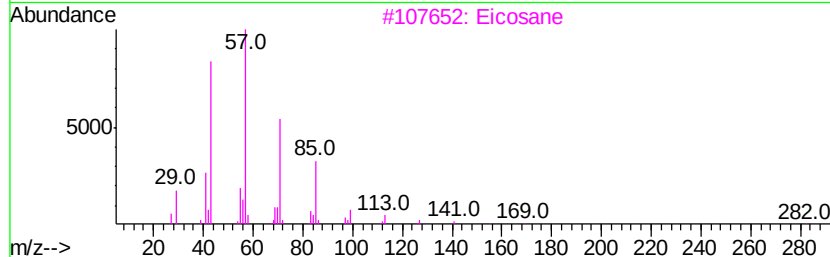
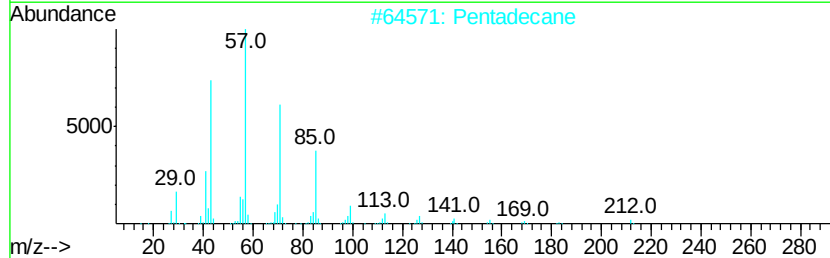
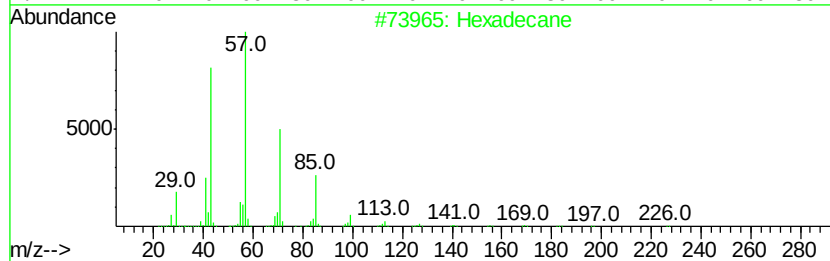
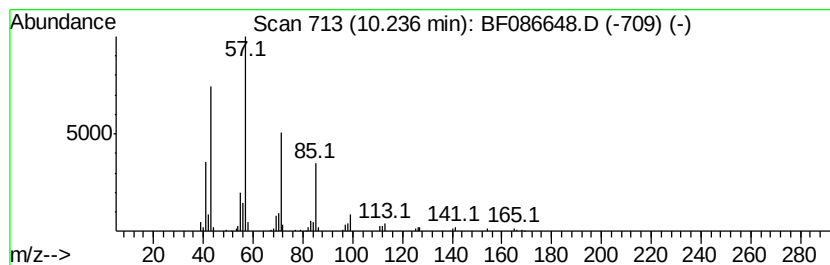
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ClientSampleId :
26WARD-2B-CS-6(10-20FTBGS)

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 6 Hexadecane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD		R.T.
10.24	2.88 ng	246147	Acenaphthene-d10		9.79
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane	226	C16H34	000544-76-3	91
2	Pentadecane	212	C15H32	000629-62-9	90
3	Eicosane	282	C20H42	000112-95-8	87
4	Tridecane	184	C13H28	000629-50-5	83
5	Tetradecane	198	C14H30	000629-59-4	80



Data Path : Z:\HPCHEM1\BNA F\DATA\BF050316\
Data File : BF086648.D
Acq On : 3 May 2016 16:57
Operator : UM/SJ
Sample : H2685-06
Misc :
ALS Vial : 38 Sample Multiplier: 1

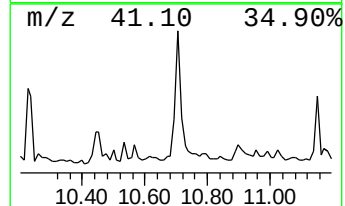
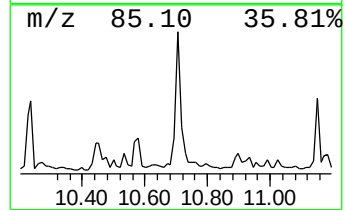
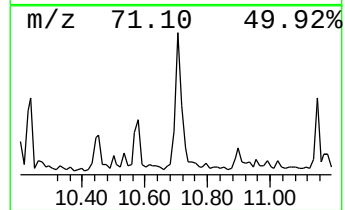
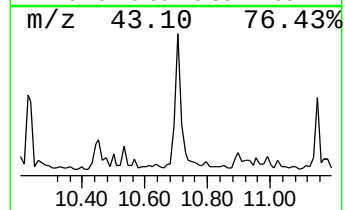
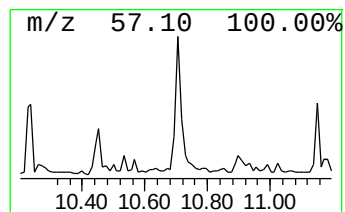
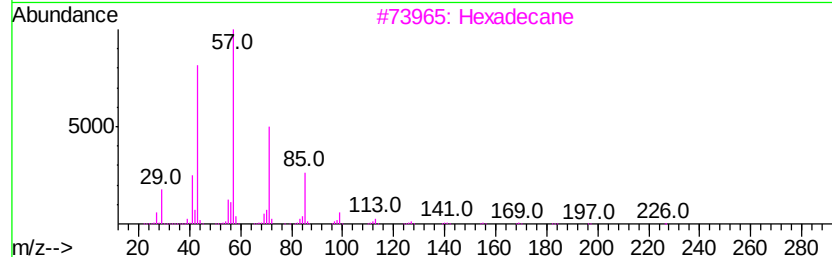
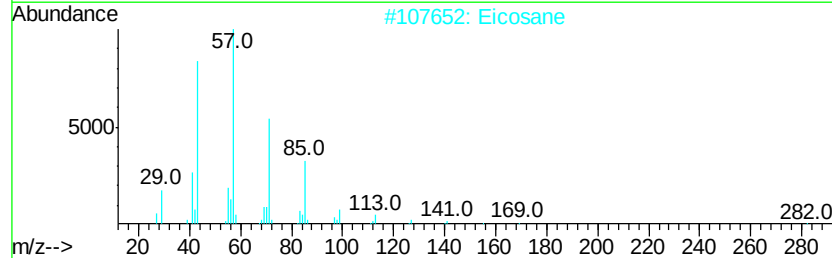
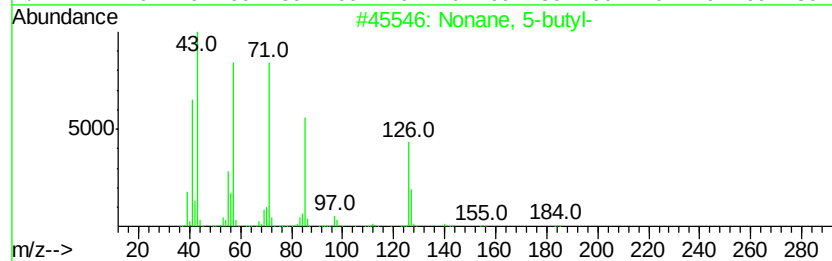
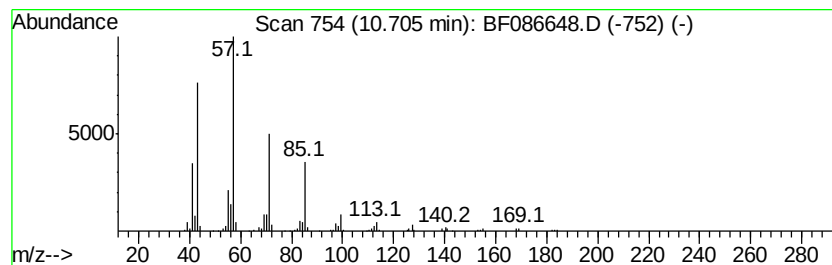
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF042716.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Nonane, 5-butyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.70	4.77 ng	374663	Phenanthrene-d10	11.26
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Nonane, 5-butyl-	184 C13H28	017312-63-9	93
2	Eicosane	282 C20H42	000112-95-8	91
3	Hexadecane	226 C16H34	000544-76-3	91
4	Dodecane, 2-methyl-6-propyl-	226 C16H34	055045-08-4	86
5	Tridecane, 6-propyl-	226 C16H34	055045-10-8	86



Data Path : Z:\HPCHEM1\BNA F\DATA\BF050316\
Data File : BF086648.D
Acq On : 3 May 2016 16:57
Operator : UM/SJ
Sample : H2685-06
Misc :
ALS Vial : 38 Sample Multiplier: 1

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ClientSampleId :
26WARD-2B-CS-6(10-20FTBGS)

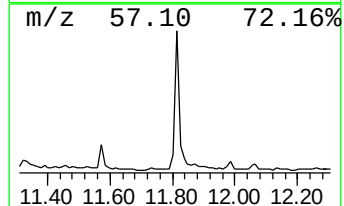
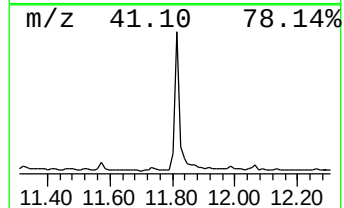
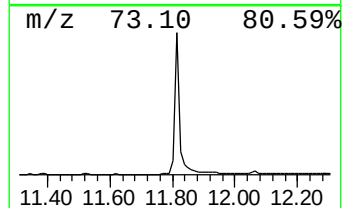
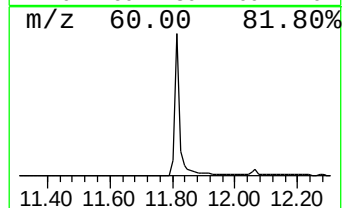
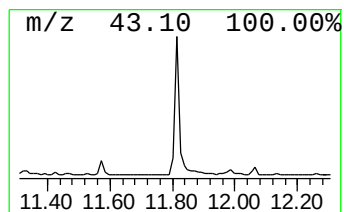
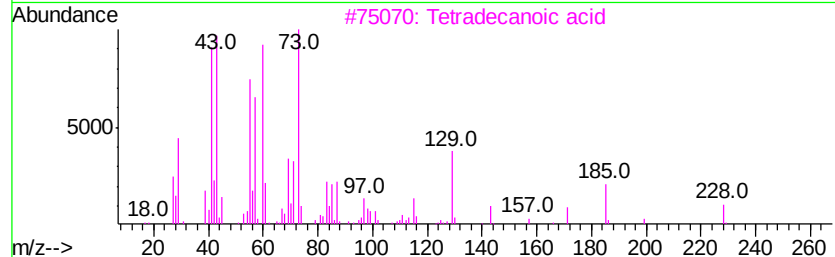
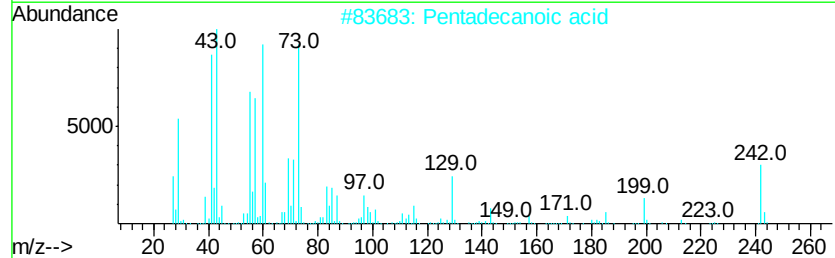
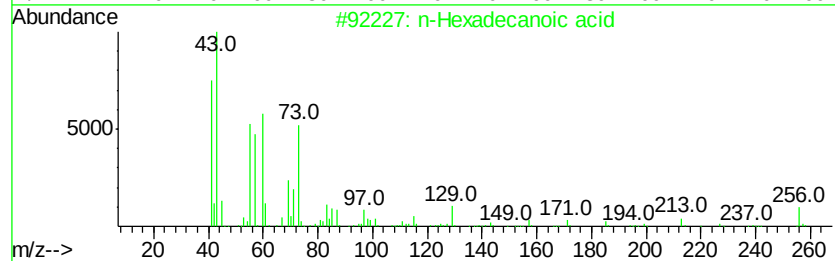
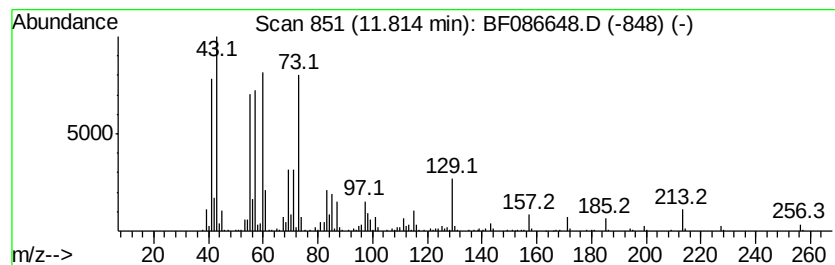
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.81	8.72 ng	685024	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	91
3	Tetradecanoic acid	228	C14H28O2	000544-63-8	83
4	n-Decanoic acid	172	C10H20O2	000334-48-5	68
5	Undecanoic acid	186	C11H22O2	000112-37-8	64



Data Path : Z:\HPCHEM1\BNA F\DATA\BF050316\
Data File : BF086648.D
Acq On : 3 May 2016 16:57
Operator : UM/SJ
Sample : H2685-06
Misc :
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Instrument :
BNA_F
ClientSampleId :
26WARD-2B-CS-6(10-20FTBGS)

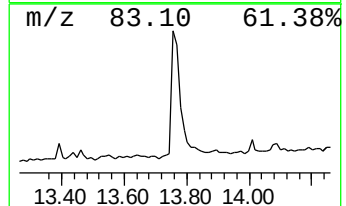
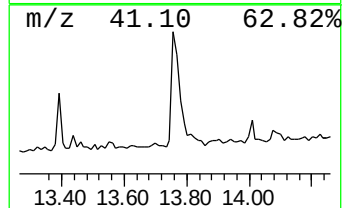
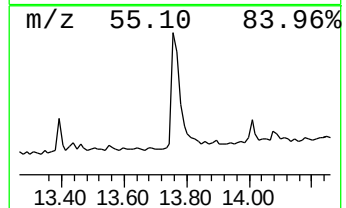
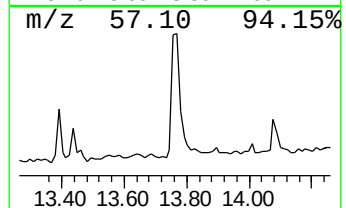
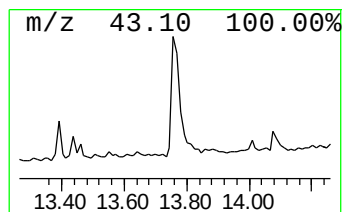
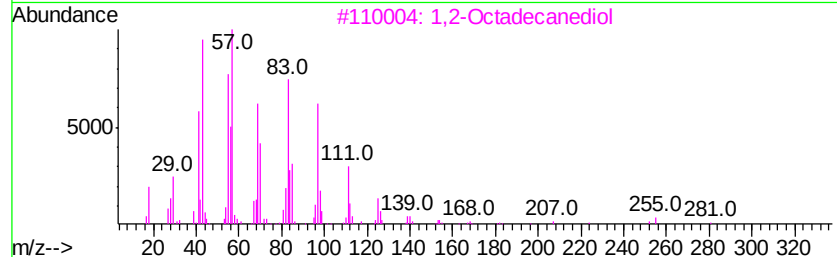
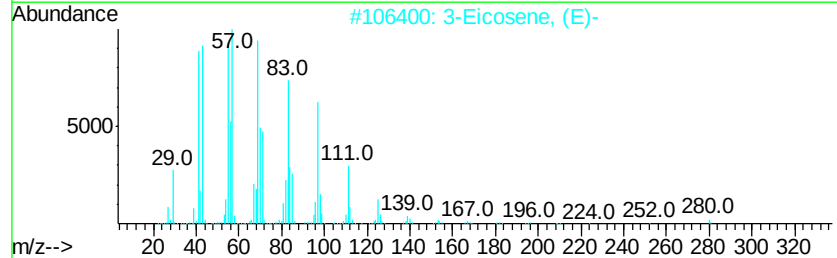
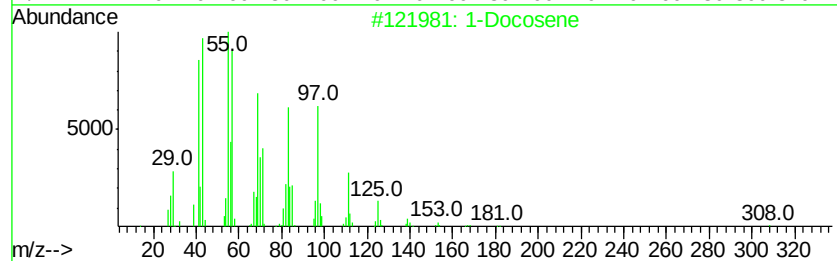
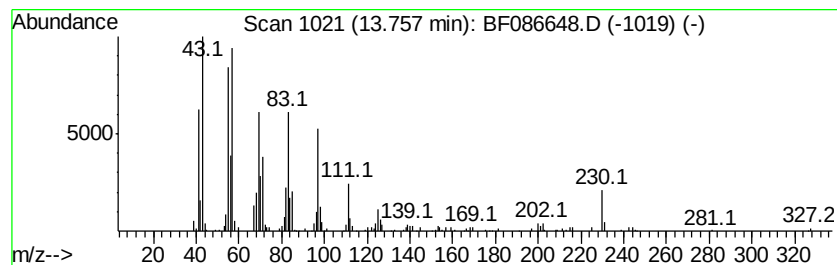
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 1-Docosene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.76	3.49 ng	307574	Chrysene-d12	13.91

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Docosene	308	C22H44	001599-67-3	93
2		3-Eicosene, (E)-	280	C20H40	074685-33-9	90
3		1,2-Octadecanediol	286	C18H38O2	020294-76-2	90
4		Cyclohexadecane	224	C16H32	000295-65-8	87
5		1-Nonadecene	266	C19H38	018435-45-5	87



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF050316\
Data File : BF086648.D
Acq On : 3 May 2016 16:57
Operator : UM/SJ
Sample : H2685-06
Misc :
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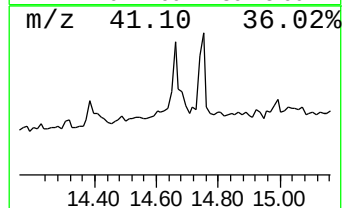
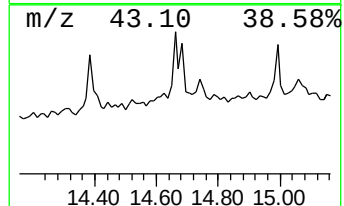
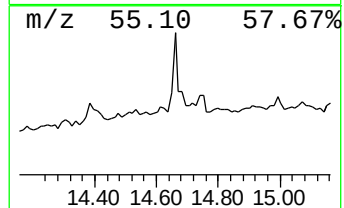
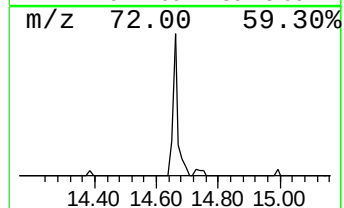
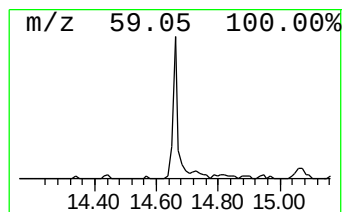
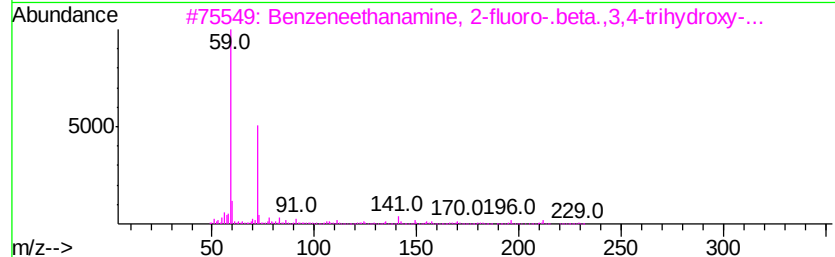
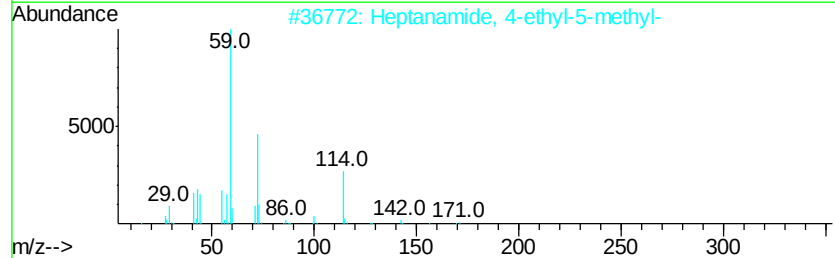
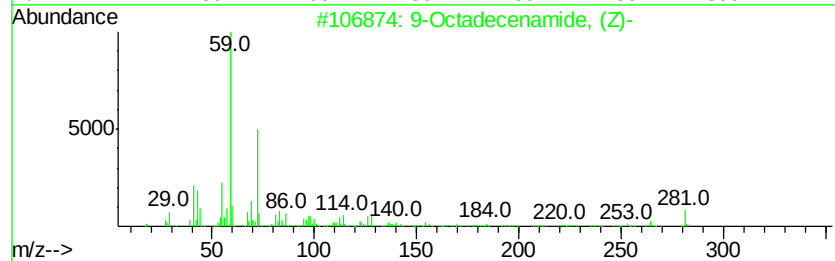
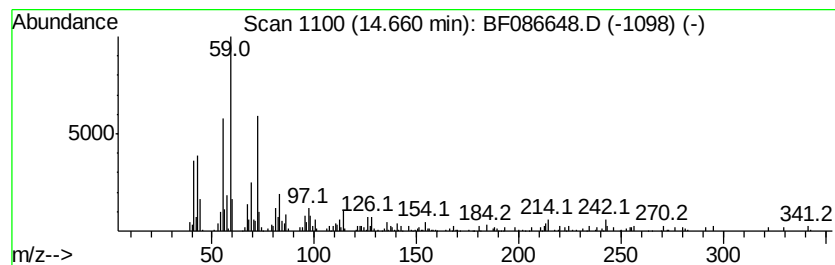
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ClientSampleId :
26WARD-2B-CS-6(10-20FTBGS)

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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 9-Octadecenamide, (Z)- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.66	2.86 ng	142678	Perylene-d12	15.30	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	93
2	Heptanamide, 4-ethyl-5-methyl-	171	C10H21NO	054789-40-1	64
3	Benzeneethanamine, 2-fluoro-.bet...	229	C11H16FN03	061338-98-5	64
4	Hexanamide	115	C6H13NO	000628-02-4	50
5	3-Cyclohexylpropionamide	155	C9H17NO	004361-29-9	47



Data Path : Z:\HPCHEM1\BNA F\DATA\BF050316\
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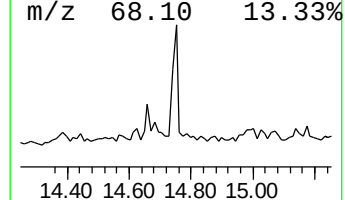
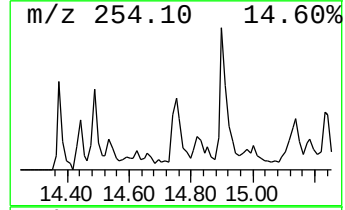
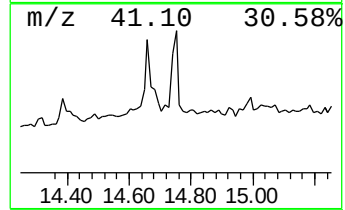
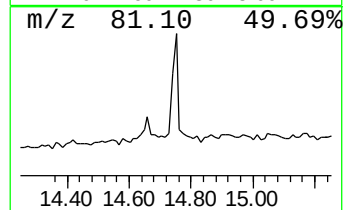
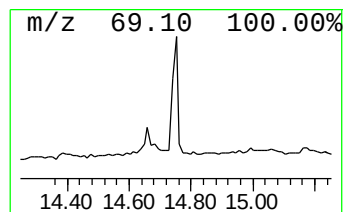
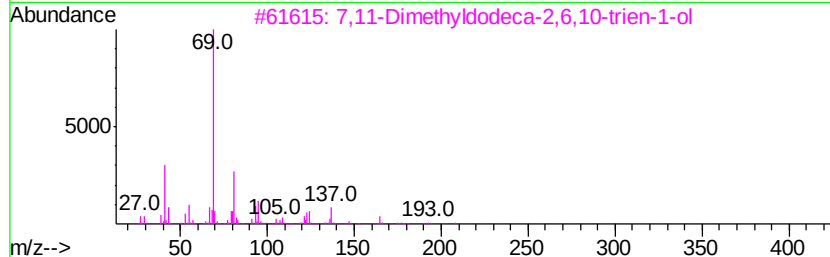
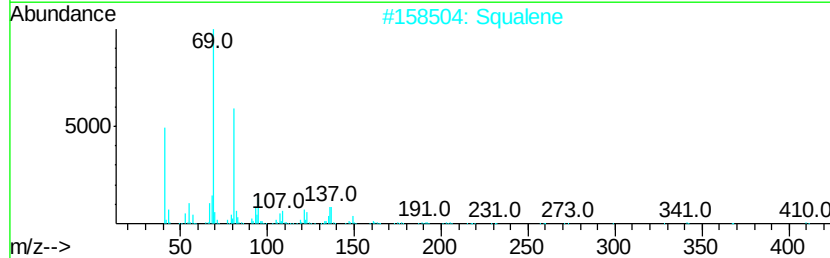
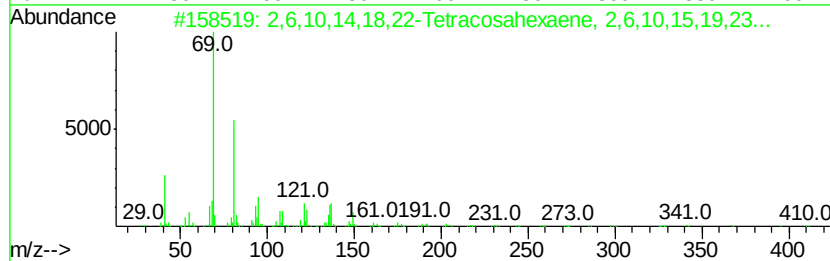
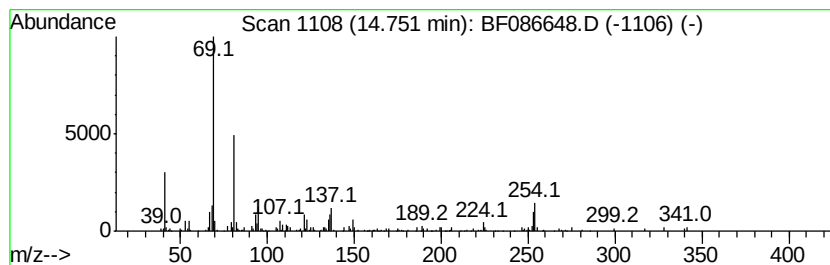
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 2,6,10,14,18,22-Tetracosahexaene... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.75	2.72 ng	135590	Perylene-d12	15.30	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,6,10,14,18,22-Tetracosahexaene...	410	C30H50	000111-02-4	68
2	Squalene	410	C30H50	007683-64-9	64
3	7,11-Dimethyldodeca-2,6,10-trien...	208	C14H24O	125529-10-4	59
4	Docosa-2,6,10,14,18-pentaen-22-a...	384	C27H44O	1000163-04-7	59
5	Benzaldehyde, 4-methoxy-3-(3,7,1...	340	C23H32O2	1000196-52-1	53



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF050316\
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 Misc :
 ALS Vial : 38 Sample Multiplier: 1

Instrument :
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 26WARD-2B-CS-6(10-20FTBGS)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF042716.M
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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	78.7	ng	5197050	1	6.75	1320750	20.0
Cyclohexasiloxane...	8.58	3.6	ng	286697	2	8.04	1604220	20.0
unknown8.90	8.90	2.8	ng	221086	2	8.04	1604220	20.0
Hexadecane	10.24	2.9	ng	246147	3	9.79	1712110	20.0
Nonane, 5-butyl-	10.70	4.8	ng	374663	4	11.26	1571400	20.0
n-Hexadecanoic acid	11.81	8.7	ng	685024	4	11.26	1571400	20.0
1-Docosene	13.76	3.5	ng	307574	5	13.91	1764510	20.0
9-Octadecenamide, ...	14.66	2.9	ng	142678	6	15.30	996100	20.0
2,6,10,14,18,22-T...	14.75	2.7	ng	135590	6	15.30	996100	20.0