

Data Path : Z:\HPCHEM1\BNA F\DATA\BF042216\
 Data File : BF086395.D
 Acq On : 23 Apr 2016 17:29
 Operator : UM/SJ
 Sample : H2640-01
 Misc :
 ALS Vial : 39 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WPG-B11(0-4)-C

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-BF042216.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	4.350	196	198	203	rBV	157833	238093	5.38%	0.528%
2	5.013	253	256	259	rBV	1007145	1466047	33.10%	3.251%
3	5.436	290	293	296	rBV	2494653	2768669	62.52%	6.140%
4	5.847	327	329	332	rBV	71065	65052	1.47%	0.144%
5	6.476	381	384	393	rBV	3213254	3866431	87.30%	8.575%
6	6.613	393	396	398	rBV	3343429	3853492	87.01%	8.546%
7	6.842	414	416	425	rBV	1189740	1196253	27.01%	2.653%
8	7.002	427	430	432	rBV	3107339	3333559	75.27%	7.393%
9	7.402	463	465	468	rBV	1693162	2393779	54.05%	5.309%
10	8.122	526	528	533	rBV	1114121	1508936	34.07%	3.346%
11	8.842	587	591	593	rVV2	45513	57650	1.30%	0.128%
12	8.933	597	599	604	rVV	45468	60824	1.37%	0.135%
13	9.208	620	623	625	rBV	3643387	4428761	100.00%	9.822%
14	9.288	628	630	632	rVB2	68640	89393	2.02%	0.198%
15	9.470	643	646	649	rBV	30305	61332	1.38%	0.136%
16	9.539	649	652	655	rVV2	67023	118646	2.68%	0.263%
17	9.596	655	657	659	rVV	615367	580486	13.11%	1.287%
18	9.733	668	669	673	rBV	39114	61252	1.38%	0.136%
19	9.813	675	676	679	rBV	116276	92396	2.09%	0.205%
20	9.882	679	682	683	rVV	1294436	1462226	33.02%	3.243%
21	9.905	683	684	686	rVV	120740	140802	3.18%	0.312%
22	10.042	693	696	698	rBV3	25696	61796	1.40%	0.137%
23	10.076	698	699	702	rVV	76296	105856	2.39%	0.235%
24	10.133	702	704	706	rVB3	36060	55927	1.26%	0.124%
25	10.282	715	717	718	rBV	49759	59434	1.34%	0.132%
26	10.316	718	720	721	rVB	162381	151923	3.43%	0.337%
27	10.419	728	729	733	rBV	88353	97848	2.21%	0.217%
28	10.533	736	739	742	rBV2	86899	133067	3.00%	0.295%
29	10.579	742	743	745	rVB	41564	47482	1.07%	0.105%
30	10.659	748	750	753	rBV	781001	857483	19.36%	1.902%
31	10.785	759	761	765	rVB	253342	315170	7.12%	0.699%
32	10.876	765	769	771	rVB4	21825	50125	1.13%	0.111%
33	10.979	775	778	783	rBV6	40396	123108	2.78%	0.273%
34	11.105	787	789	793	rVB5	20935	47170	1.07%	0.105%

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Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF042216.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	11.231	796	800	801 rBV	164021	182161	4.11%	0.404%
36	11.265	801	803	806 rVB2	102839	144548	3.26%	0.321%
37	11.356	809	811	816 rBV2	1073934	1967142	44.42%	4.363%
38	11.436	816	818	823 rVB	158368	243103	5.49%	0.539%
39	11.608	830	833	836 rBV2	40430	85528	1.93%	0.190%
40	11.654	836	837	839 rVV	74702	83658	1.89%	0.186%
41	11.859	853	855	856 rVV	103052	100218	2.26%	0.222%
42	11.882	856	857	859 rVV	84917	113067	2.55%	0.251%
43	11.928	859	861	863 rVV2	44796	64352	1.45%	0.143%
44	11.974	863	865	869 rVB2	145218	265496	5.99%	0.589%
45	12.065	872	873	874 rBV	86626	64238	1.45%	0.142%
46	12.156	879	881	882 rBV2	41775	57687	1.30%	0.128%
47	12.351	895	898	901 rBV3	37432	85528	1.93%	0.190%
48	12.408	901	903	904 rBV	85232	86444	1.95%	0.192%
49	12.454	904	907	909 rVB3	147764	259944	5.87%	0.576%
50	12.488	909	910	913 rBV2	37122	51936	1.17%	0.115%
51	12.568	914	917	920 rBV	807308	821892	18.56%	1.823%
52	12.797	934	937	938 rBV	707091	815172	18.41%	1.808%
53	12.819	938	939	946 rVB3	155161	260729	5.89%	0.578%
54	12.945	948	950	954 rBV	2953680	3220348	72.71%	7.142%
55	13.128	964	966	968 rVB	102402	120598	2.72%	0.267%
56	13.185	968	971	973 rBV2	228816	317615	7.17%	0.704%
57	13.231	973	975	979 rVV3	73681	149119	3.37%	0.331%
58	13.528	998	1001	1003 rBV	203163	341161	7.70%	0.757%
59	13.848	1027	1029	1034 rVB	436108	594386	13.42%	1.318%
60	13.985	1038	1041	1048 rVV2	831135	1831307	41.35%	4.061%
61	14.168	1055	1057	1059 rBV	318519	305418	6.90%	0.677%
62	14.465	1081	1083	1085 rBV	234586	335119	7.57%	0.743%
63	14.774	1107	1110	1113 rBV2	216748	368906	8.33%	0.818%
64	15.014	1128	1131	1135 rBV	191456	449389	10.15%	0.997%
65	15.094	1136	1138	1141 rVB2	221095	282561	6.38%	0.627%
66	15.414	1164	1166	1168 rBV	445385	616812	13.93%	1.368%
67	16.077	1222	1224	1233 rVB2	181104	455960	10.30%	1.011%

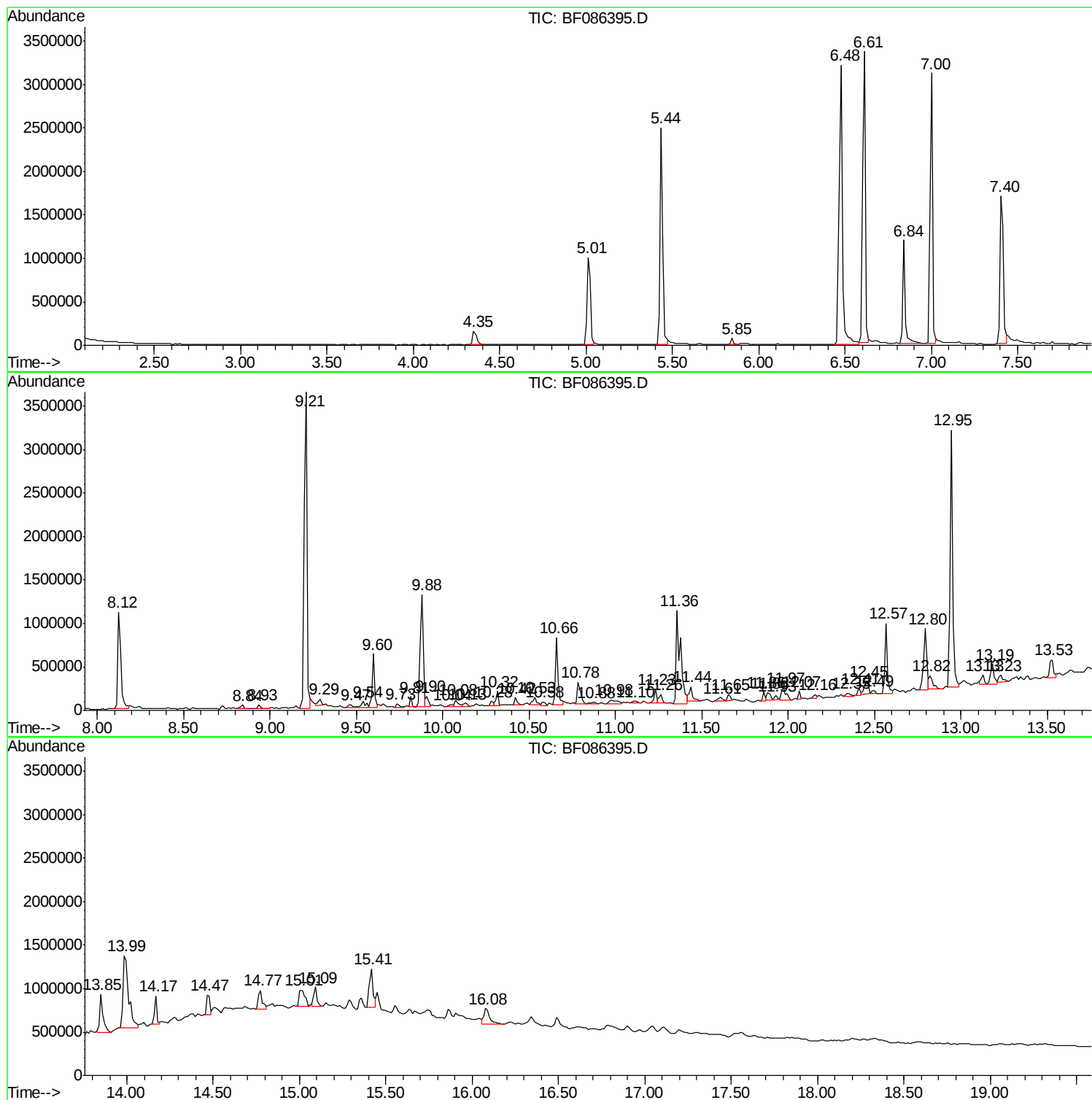
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Instrument :
BNA_F
ClientSampled :
WPG-B11(0-4)-C

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF042216.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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Instrument :
BNA_F
ClientSampleID :
WPG-B11(0-4)-C

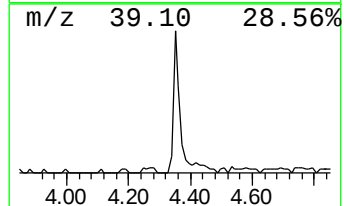
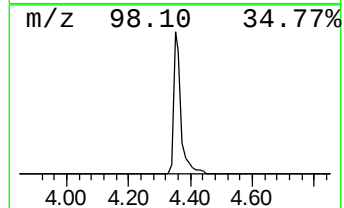
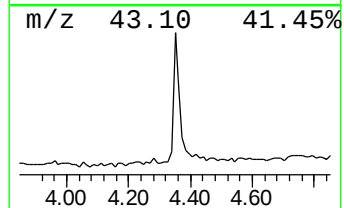
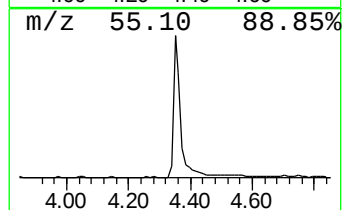
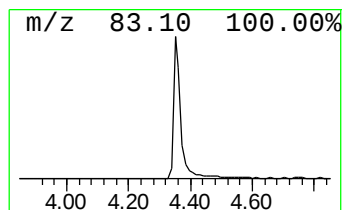
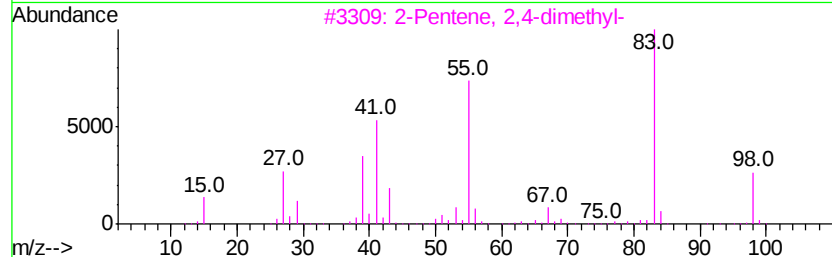
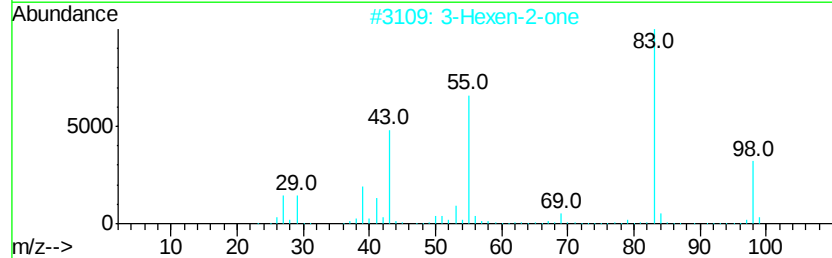
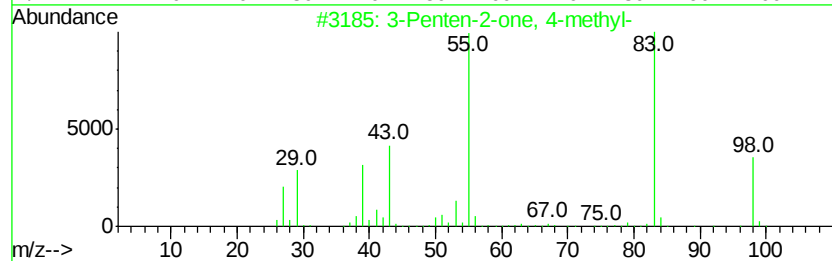
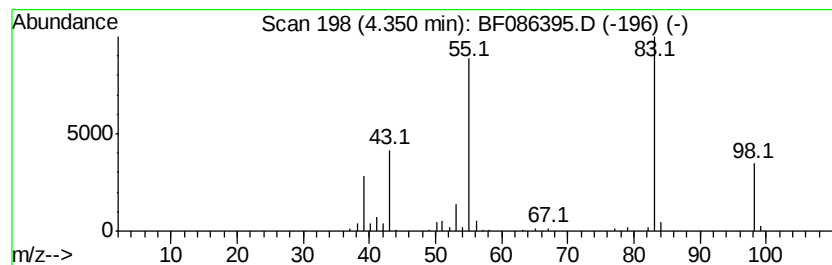
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF042216.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 3-Penten-2-one, 4-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.35	3.98 ng	238093	1,4-Dichlorobenzene-d4	6.84

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Penten-2-one, 4-methyl-	98	C6H10O	000141-79-7	91
2			3-Hexen-2-one	98	C6H10O	000763-93-9	90
3			2-Pentene, 2,4-dimethyl-	98	C7H14	000625-65-0	80
4			2-Pentene, 3,4-dimethyl-	98	C7H14	024910-63-2	80
5			3,4-Dimethyl-2-pentene(c,t)	98	C7H14	1000118-16-5	78



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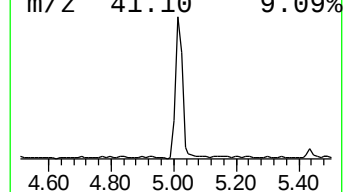
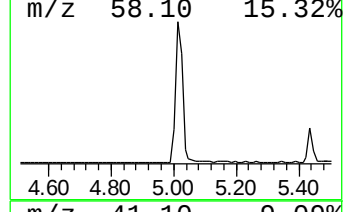
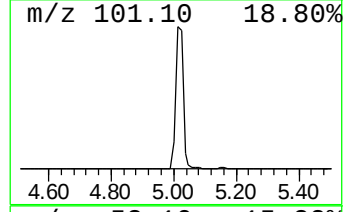
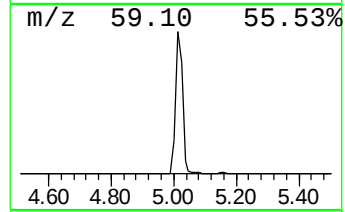
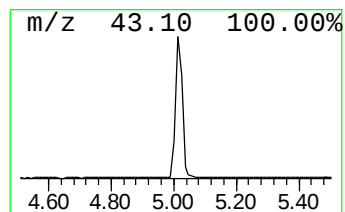
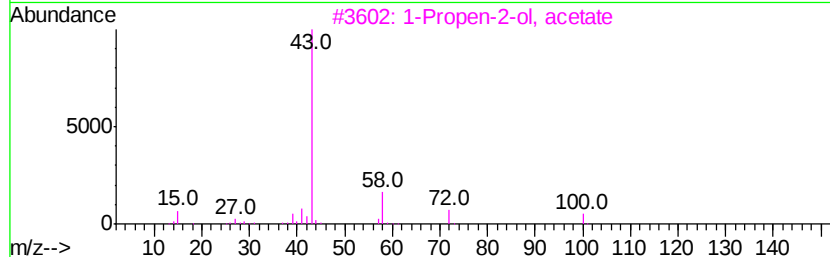
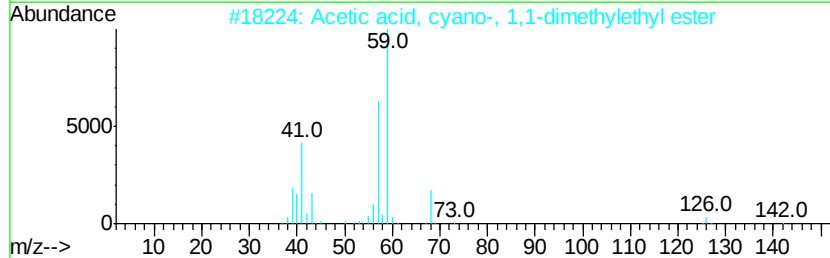
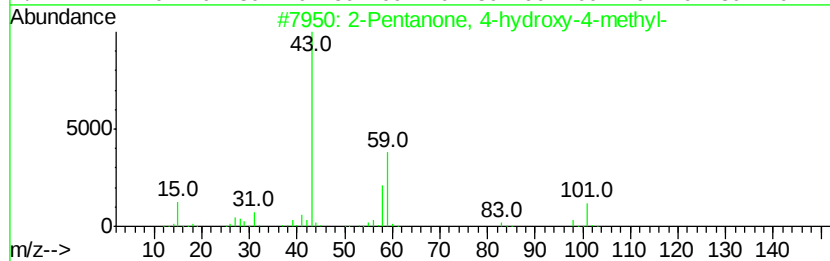
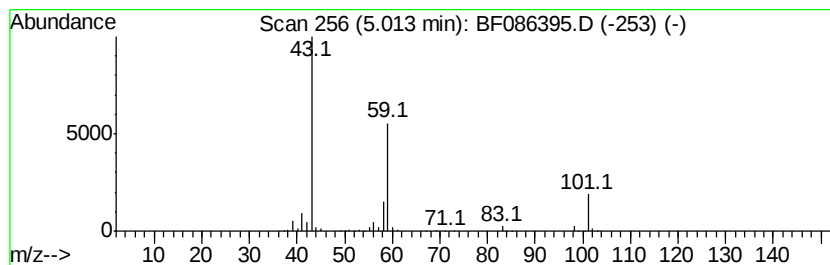
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF042216.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.01	24.51 ng	1466050	1,4-Dichlorobenzene-d4	6.84

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2		Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	23
3		1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10
4		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
5		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9



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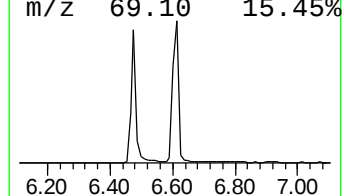
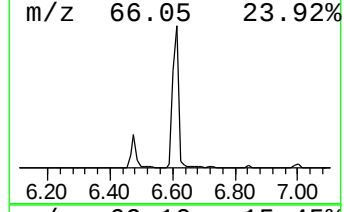
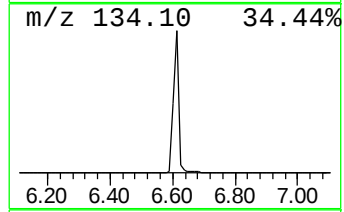
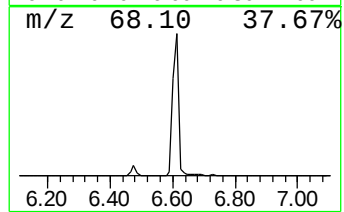
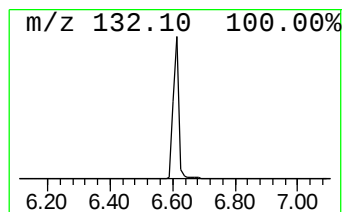
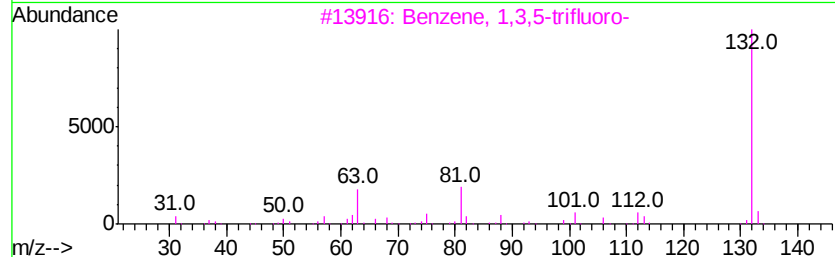
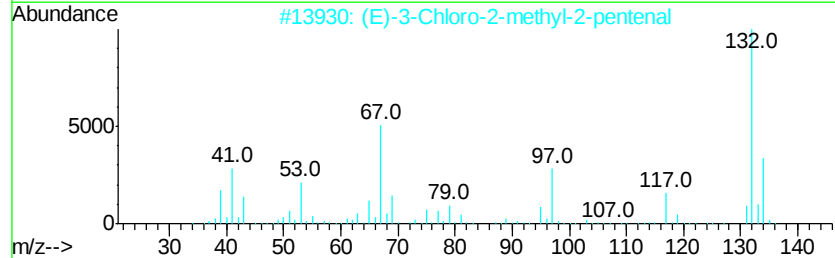
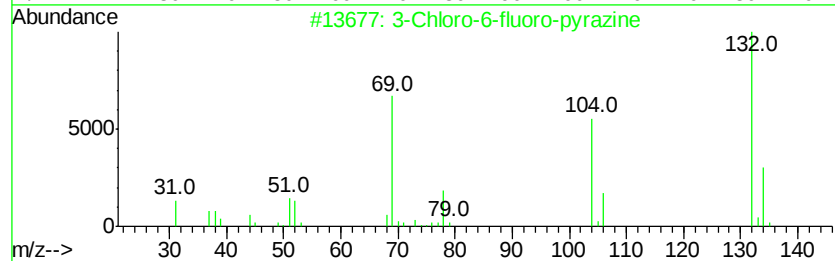
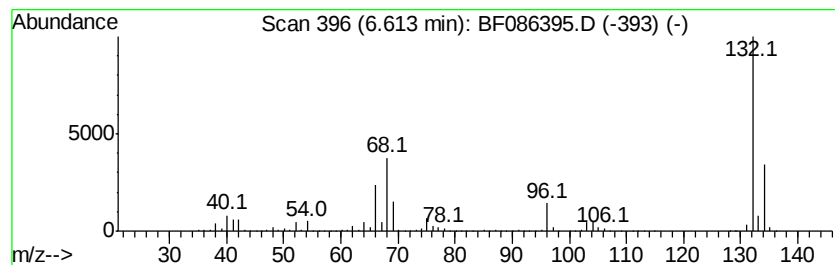
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 unknown6.61 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.61	64.43 ng	3853490	1,4-Dichlorobenzene-d4	6.84

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	35
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
3		Benzene, 1,3,5-trifluoro-	132	C6H3F3	000372-38-3	9
4		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
5		4-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-60-2	9



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ALS Vial : 39 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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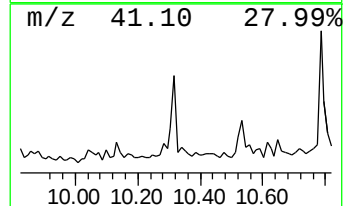
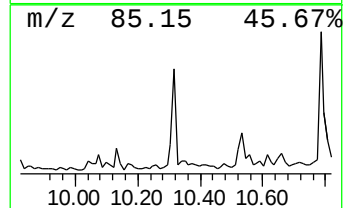
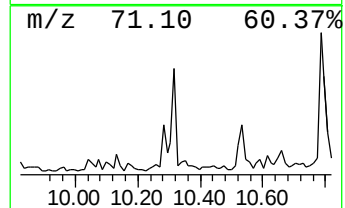
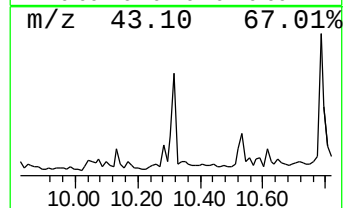
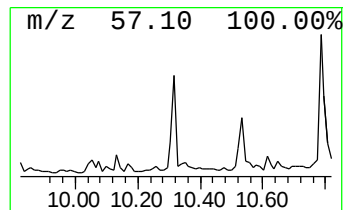
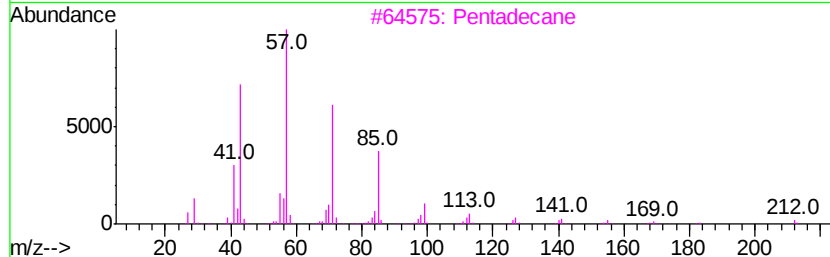
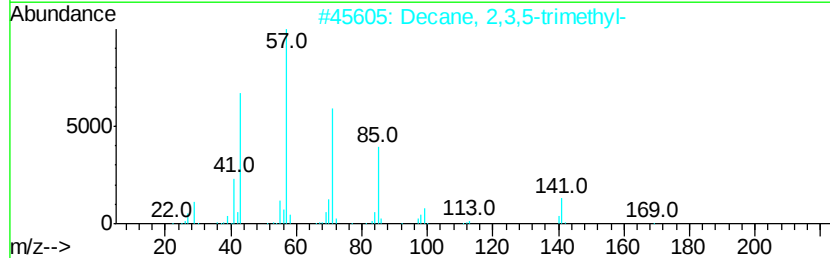
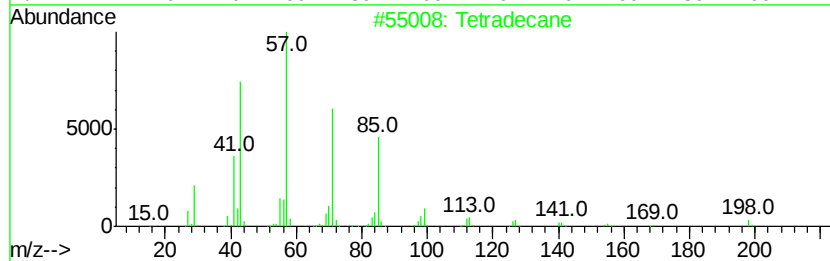
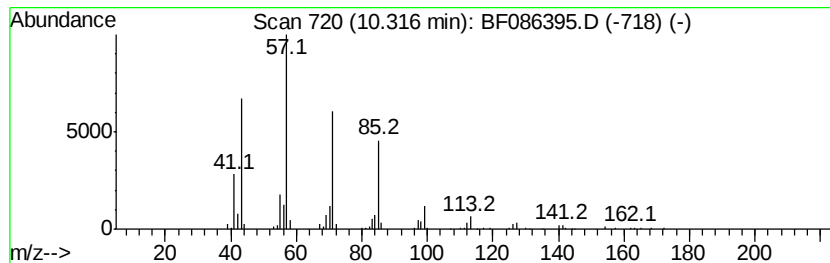
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF042216.M
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Tetradecane Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.32	2.08 ng	151923	Acenaphthene-d10	9.88

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetradecane	198	C14H30	000629-59-4	91
2		Decane, 2,3,5-trimethyl-	184	C13H28	062238-11-3	90
3		Pentadecane	212	C15H32	000629-62-9	90
4		Tridecane	184	C13H28	000629-50-5	90
5		Heptadecane	240	C17H36	000629-78-7	87



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ALS Vial : 39 Sample Multiplier: 1

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ClientSampleId :
WPG-B11(0-4)-C

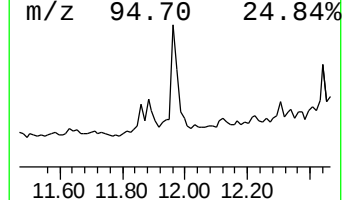
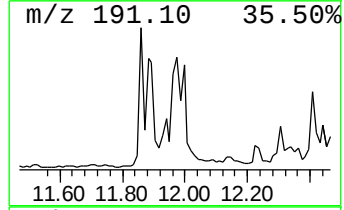
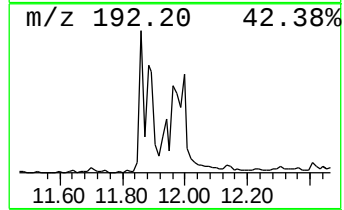
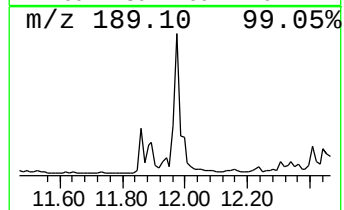
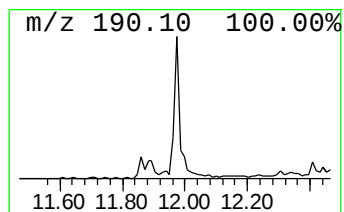
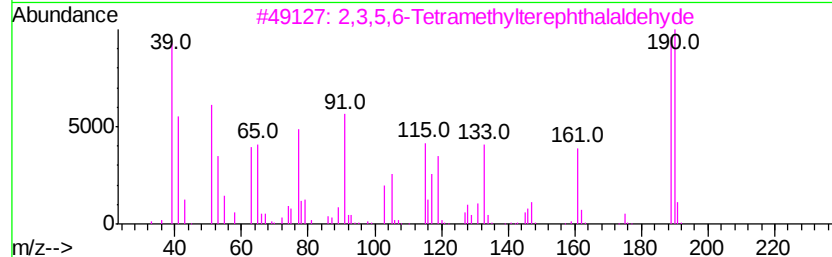
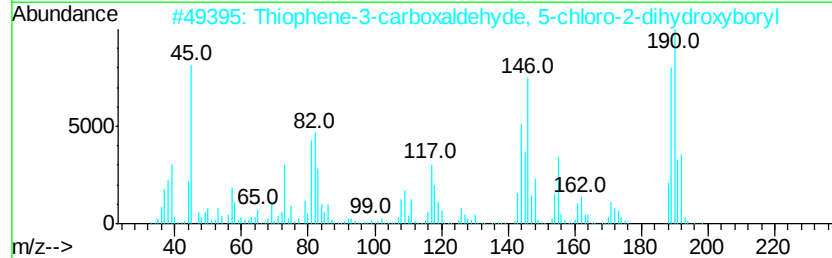
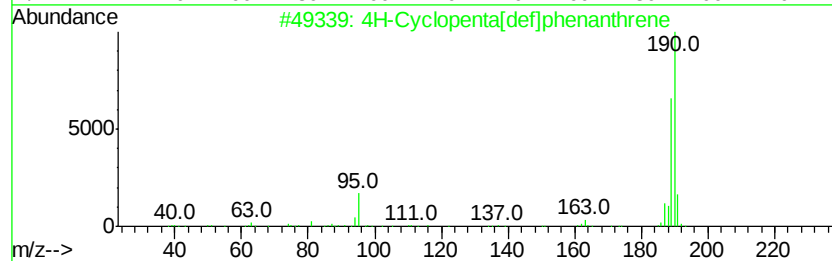
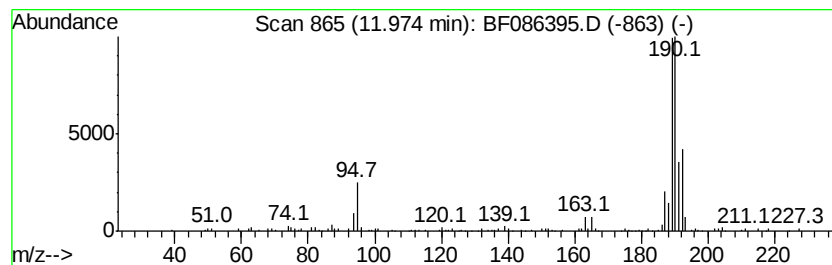
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 4H-Cyclopenta[def]phenanthrene Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.97	2.70 ng	265496	Phenanthrene-d10	11.36

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	81
2		Thiophene-3-carboxaldehyde, 5-ch...	190	C5H4BClO3S	036155-87-0	50
3		2,3,5,6-Tetramethylterephthalald...	190	C12H14O2	007072-01-7	47
4		Methyl diselenide	190	C2H6Se2	007101-31-7	43
5		Pyrazole-4-carboxaldehyde, 3-(4-...	190	C10H7FN2O	306936-57-2	40



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF042216\
Data File : BF086395.D
Acq On : 23 Apr 2016 17:29
Operator : UM/SJ
Sample : H2640-01
Misc :
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Instrument :
BNA_F
ClientSampleId :
WPG-B11(0-4)-C

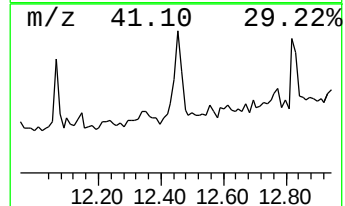
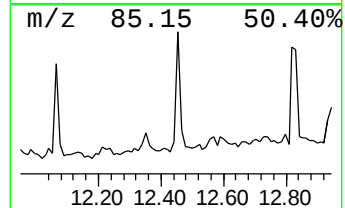
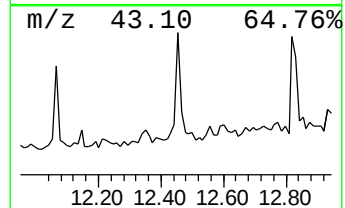
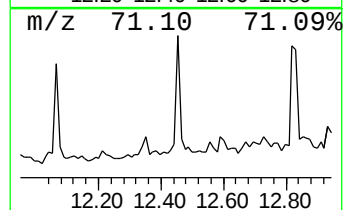
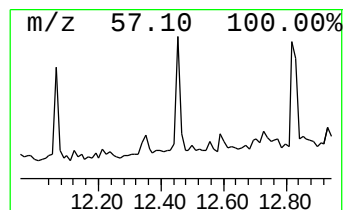
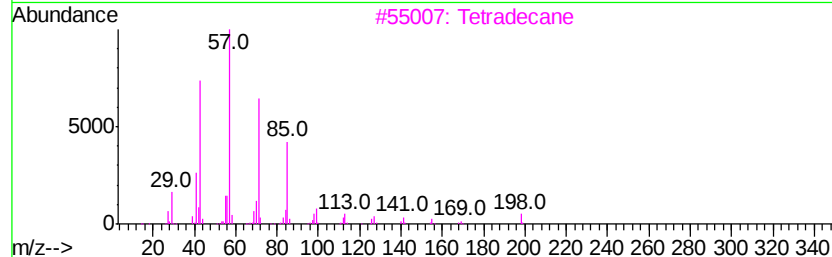
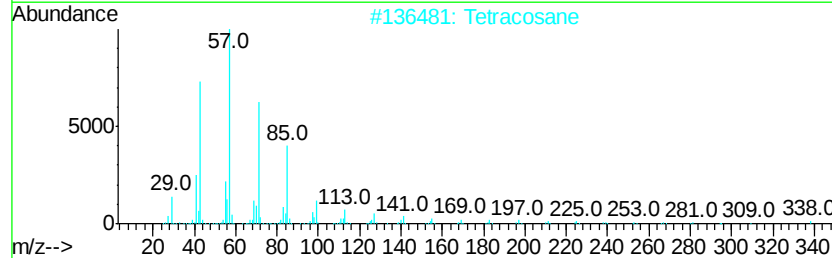
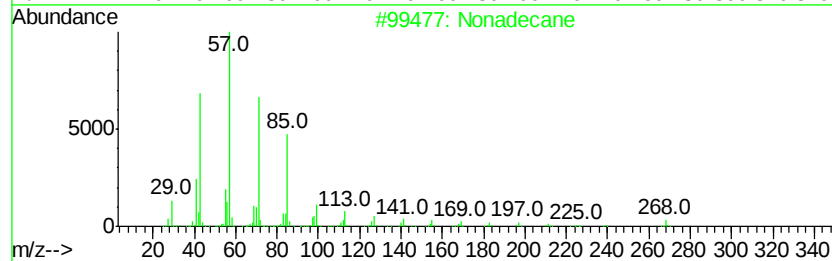
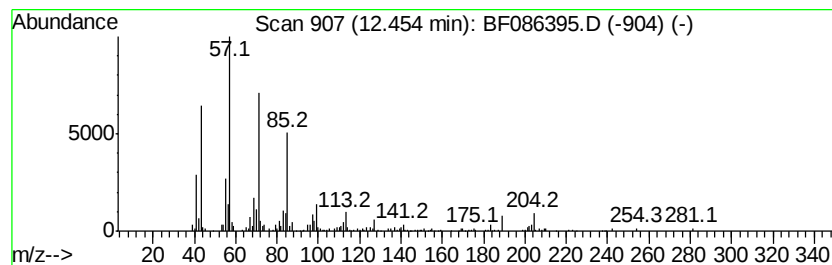
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Nonadecane Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.45	2.64 ng	259944	Phenanthrene-d10	11.36

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonadecane		268	C19H40	000629-92-5	83
2	Tetracosane		338	C24H50	000646-31-1	83
3	Tetradecane		198	C14H30	000629-59-4	83
4	Heptacosane		380	C27H56	000593-49-7	83
5	Dodecane		170	C12H26	000112-40-3	83



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF042216\
Data File : BF086395.D
Acq On : 23 Apr 2016 17:29
Operator : UM/SJ
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Misc :
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Instrument :
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WPG-B11(0-4)-C

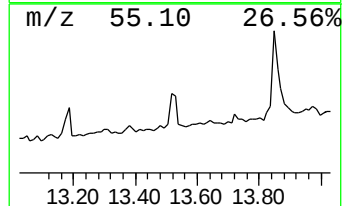
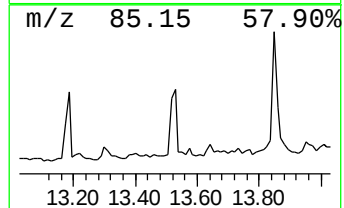
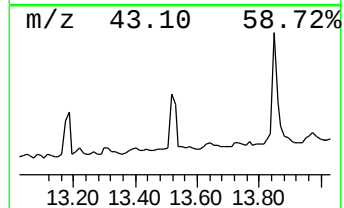
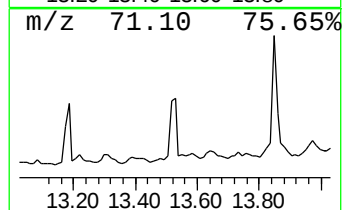
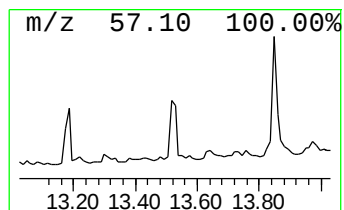
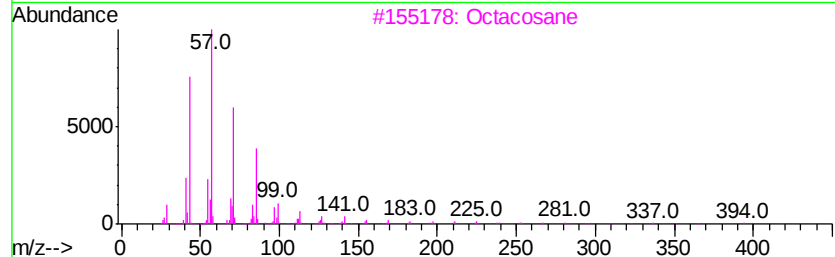
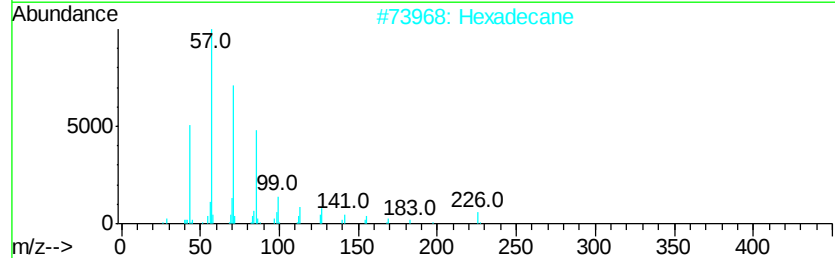
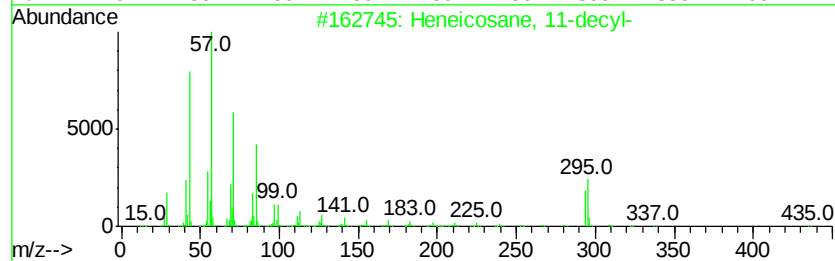
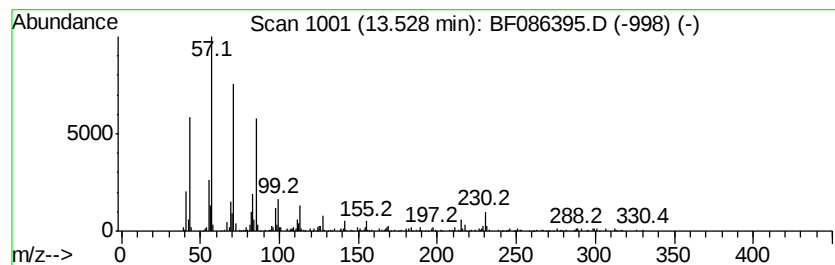
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Heneicosane, 11-decyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.53	3.73 ng	341161	Chrysene-d12	14.00

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane, 11-decyl-	437	C31H64	055320-06-4	86
2	Hexadecane	226	C16H34	000544-76-3	83
3	Octacosane	394	C28H58	000630-02-4	83
4	Tetratetracontane	619	C44H90	007098-22-8	81
5	Tridecane, 7-hexyl-	268	C19H40	007225-66-3	74



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF042216\
Data File : BF086395.D
Acq On : 23 Apr 2016 17:29
Operator : UM/SJ
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Misc :
ALS Vial : 39 Sample Multiplier: 1

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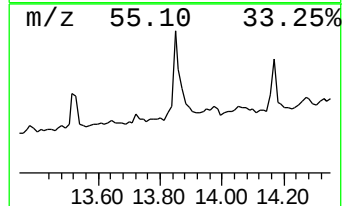
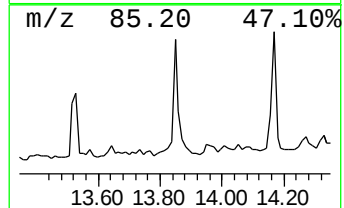
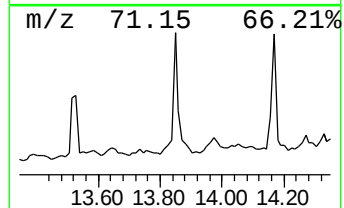
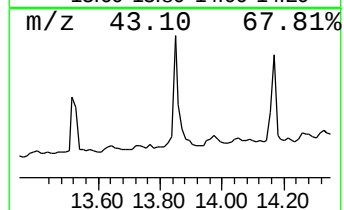
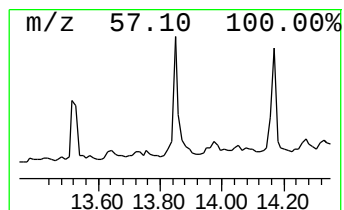
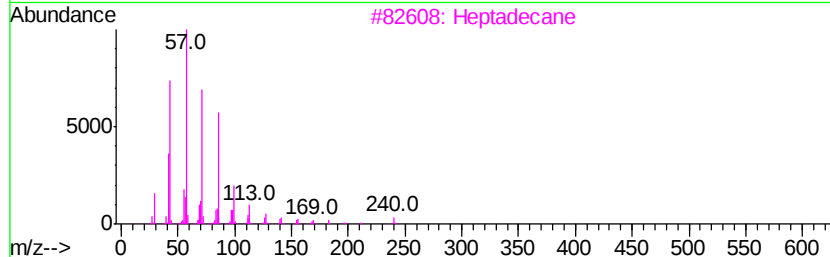
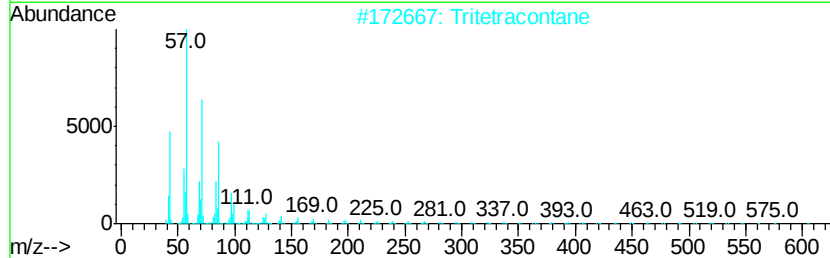
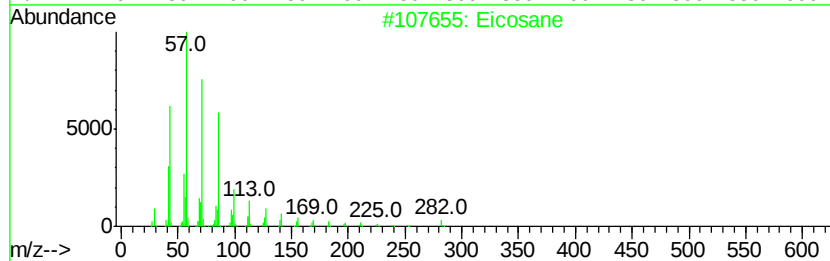
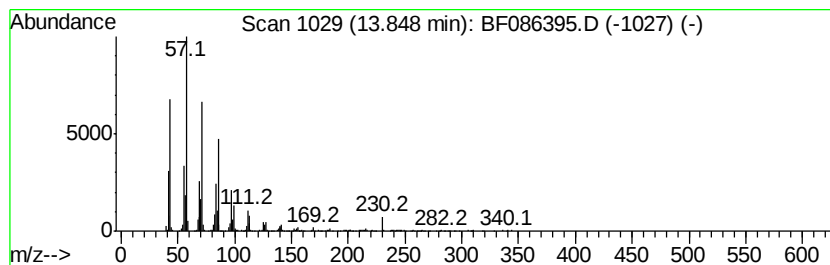
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Eicosane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.85	6.49 ng	594386	Chrysene-d12	14.00

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane			282	C20H42	000112-95-8	95
2	Tritetracontane			605	C43H88	007098-21-7	87
3	Heptadecane			240	C17H36	000629-78-7	70
4	Tridecane, 1-iodo-			310	C13H27I	035599-77-0	70
5	Octadecane			254	C18H38	000593-45-3	62



Data Path : Z:\HPCHEM1\BNA F\DATA\BF042216\
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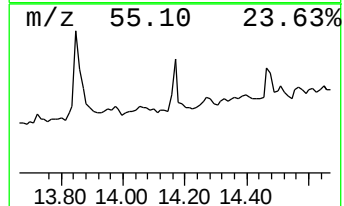
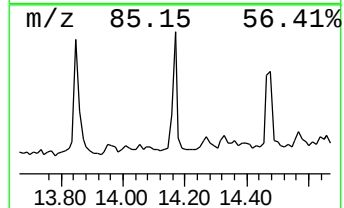
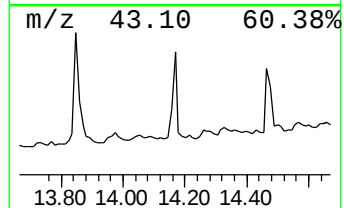
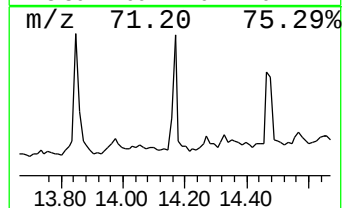
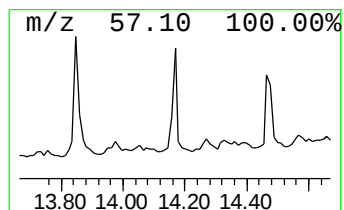
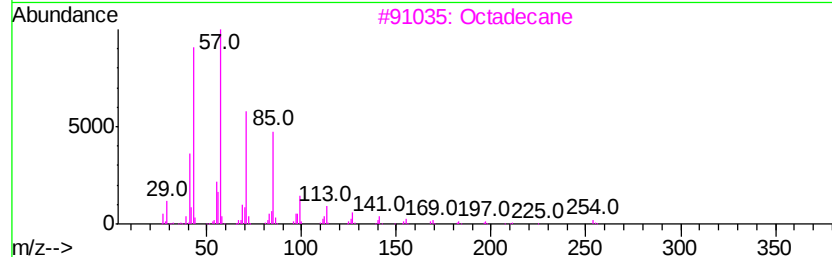
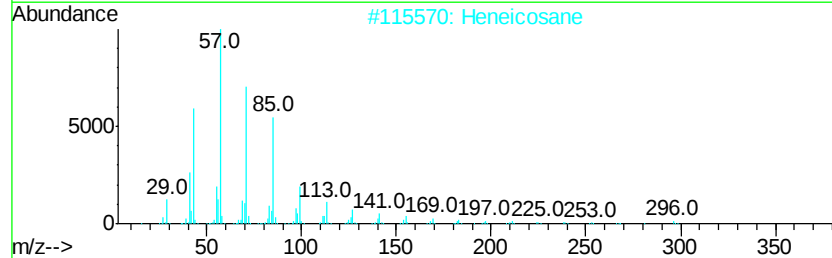
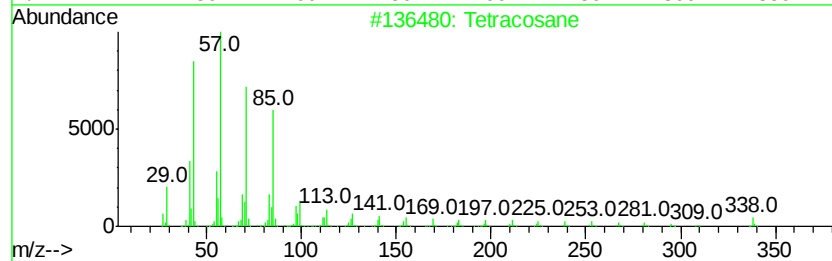
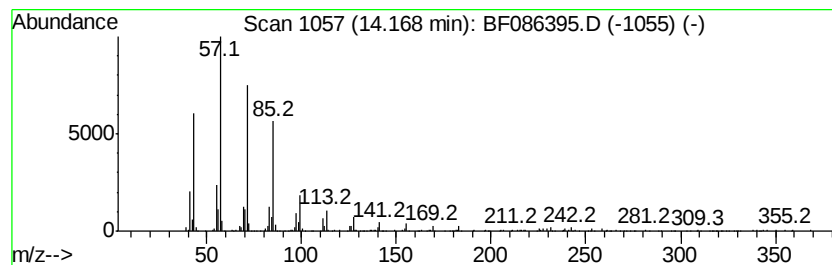
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Tetracosane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.17	3.34 ng	305418	Chrysene-d12	14.00

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetracosane	338	C24H50	000646-31-1	94
2		Heneicosane	296	C21H44	000629-94-7	91
3		Octadecane	254	C18H38	000593-45-3	91
4		Nonacosane	408	C29H60	000630-03-5	91
5		Pentadecane, 8-heptyl-	310	C22H46	071005-15-7	91



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF042216\
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Operator : UM/SJ
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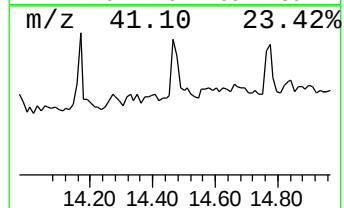
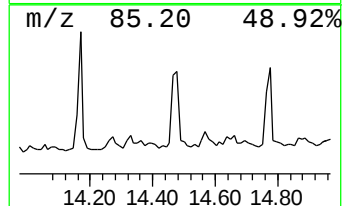
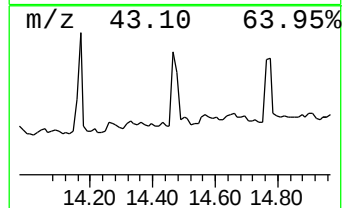
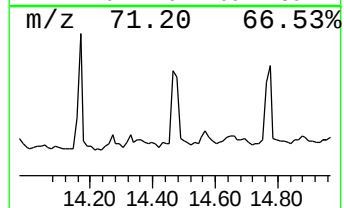
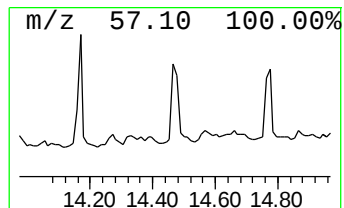
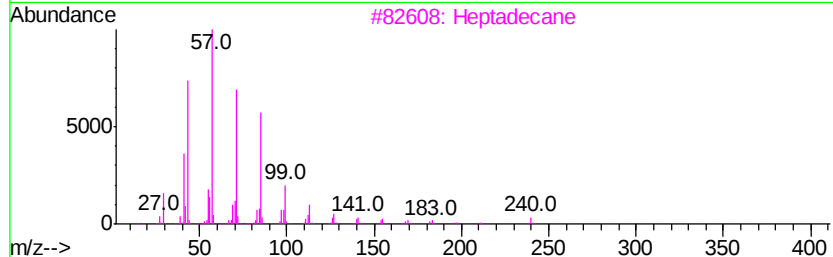
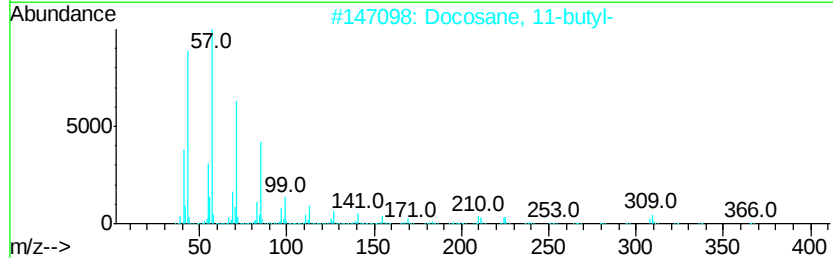
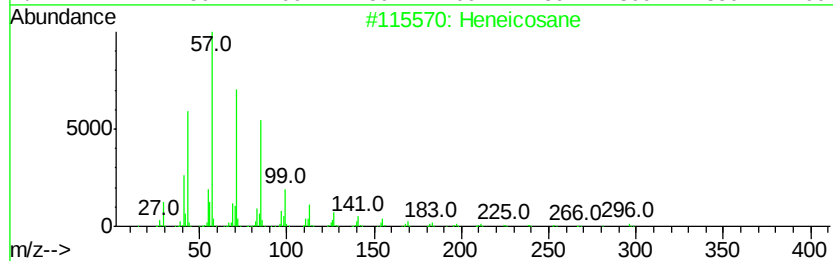
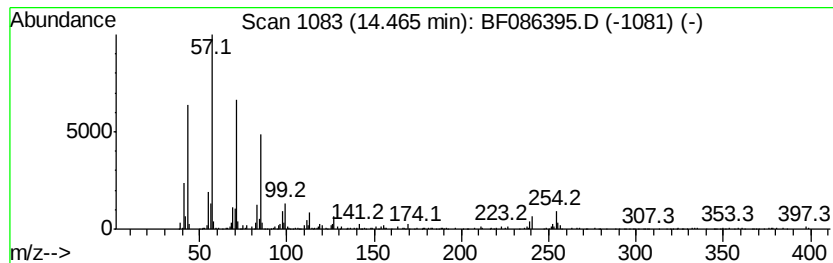
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Heneicosane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.47	3.66 ng	335119	Chrysene-d12	14.00	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane	296	C21H44	000629-94-7	95
2	Docosane, 11-butyl-	366	C26H54	013475-76-8	91
3	Heptadecane	240	C17H36	000629-78-7	91
4	Docosane, 9-butyl-	366	C26H54	055282-14-9	87
5	Octacosane	394	C28H58	000630-02-4	87



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF042216\
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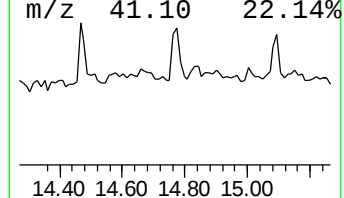
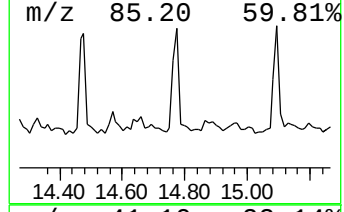
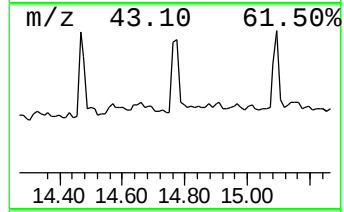
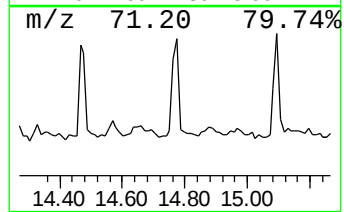
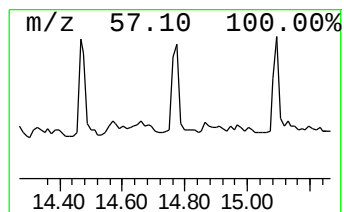
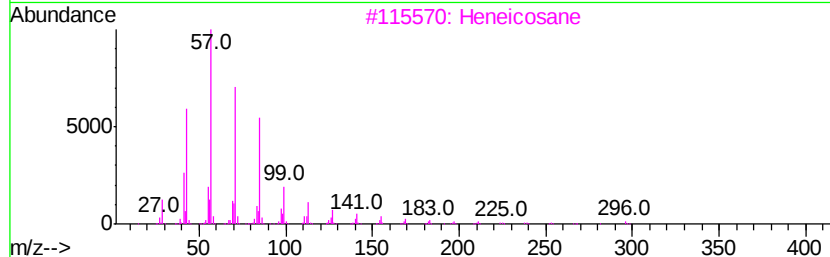
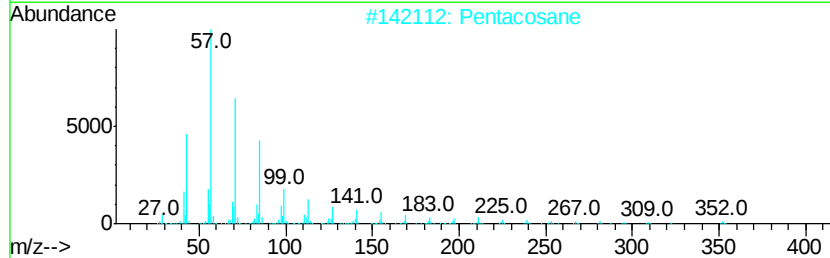
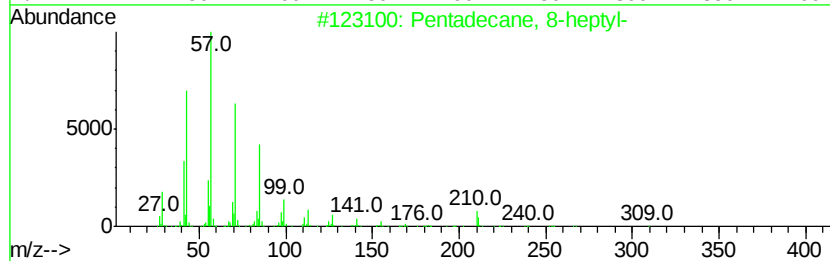
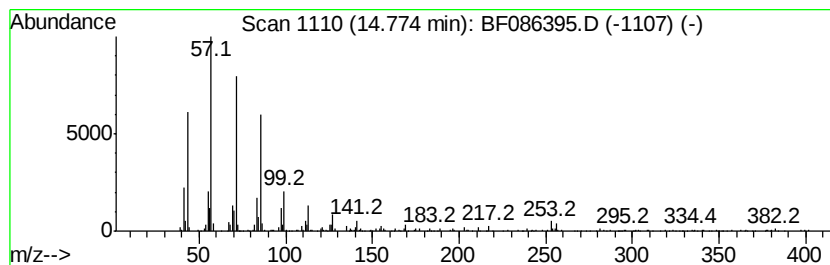
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Pentadecane, 8-heptyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.77	11.96 ng	368906	Perylene-d12	15.41

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentadecane, 8-heptyl-	310	C22H46	071005-15-7	91
2		Pentacosane	352	C25H52	000629-99-2	91
3		Heneicosane	296	C21H44	000629-94-7	91
4		Nonacosane	408	C29H60	000630-03-5	87
5		Nonadecane	268	C19H40	000629-92-5	87



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF042216\
Data File : BF086395.D
Acq On : 23 Apr 2016 17:29
Operator : UM/SJ
Sample : H2640-01
Misc :
ALS Vial : 39 Sample Multiplier: 1

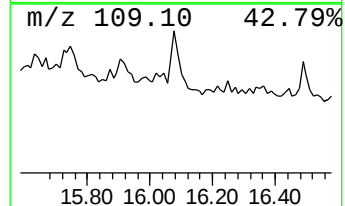
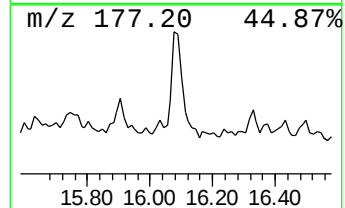
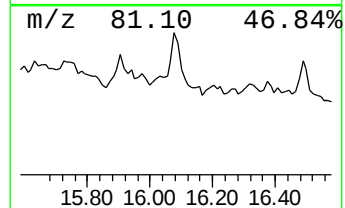
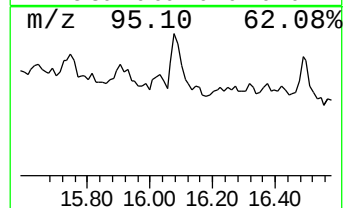
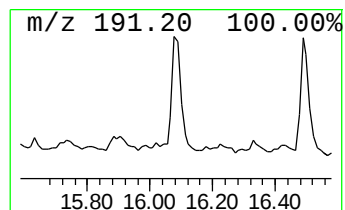
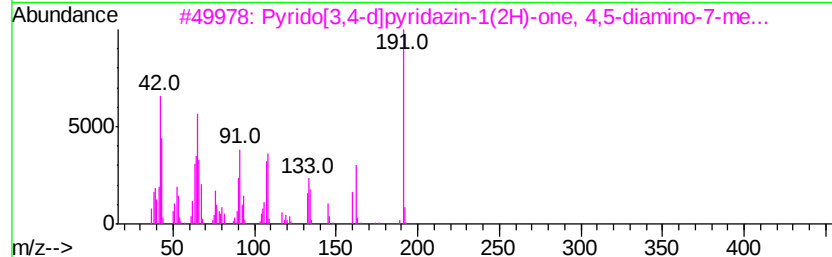
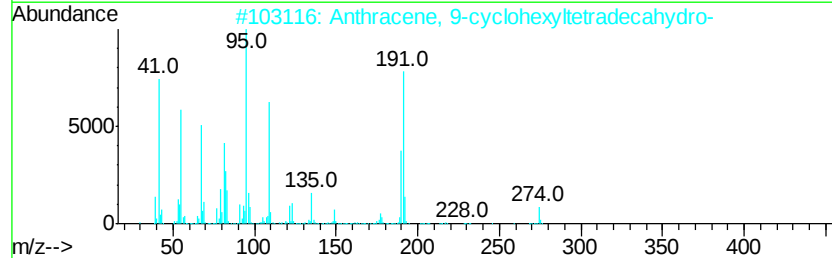
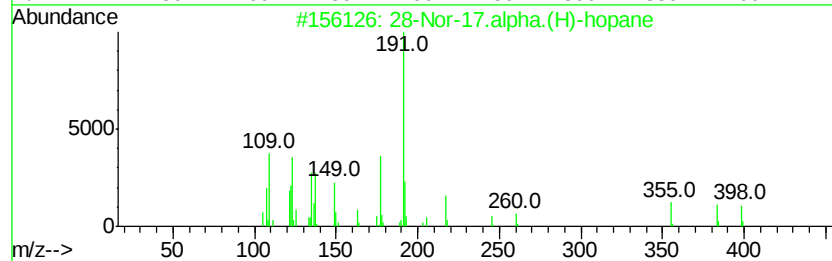
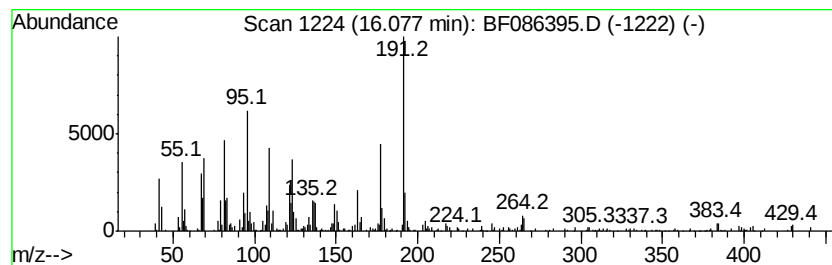
Instrument :
BNA_F
ClientSampleId :
WPG-B11(0-4)-C

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF042216.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 28-Nor-17.alpha.(H)-hopane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.08	14.78 ng	455960	Perylene-d12	15.41
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	28-Nor-17.alpha.(H)-hopane	398 C29H50	053584-60-4	72
2	Anthracene, 9-cyclohexyltetradec...	274 C20H34	055255-70-4	50
3	Pyrido[3,4-d]pyridazin-1(2H)-one...	191 C8H9N5O	1000263-69-7	35
4	5-(1-Isopropenyl-4,5-dimethylbic...	332 C22H36O2	1000195-69-8	35
5	5H-Benzo[def]carbazole	191 C14H9N	000203-65-6	27



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF042216\
Data File : BF086395.D
Acq On : 23 Apr 2016 17:29
Operator : UM/SJ
Sample : H2640-01
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
WPG-B11(0-4)-C

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF042216.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
3-Penten-2-one, 4...	4.35	4.0 ng		238093	1	6.84	1196250	20.0
2-Pentanone, 4-hy...	5.01	24.5 ng		1466050	1	6.84	1196250	20.0
unknown6.61	6.61	64.4 ng		3853490	1	6.84	1196250	20.0
Tetradecane	10.32	2.1 ng		151923	3	9.88	1462230	20.0
4H-Cyclopenta[def...	11.97	2.7 ng		265496	4	11.36	1967140	20.0
Nonadecane	12.45	2.6 ng		259944	4	11.36	1967140	20.0
Heneicosane, 11-d...	13.53	3.7 ng		341161	5	14.00	1831310	20.0
Eicosane	13.85	6.5 ng		594386	5	14.00	1831310	20.0
Tetracosane	14.17	3.3 ng		305418	5	14.00	1831310	20.0
Heneicosane	14.47	3.7 ng		335119	5	14.00	1831310	20.0
Pentadecane, 8-he...	14.77	12.0 ng		368906	6	15.41	616812	20.0
28-Nor-17.alpha.(...	16.08	14.8 ng		455960	6	15.41	616812	20.0