

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF020116\
 Data File : BF084715.D
 Acq On : 1 Feb 2016 18:00
 Operator : SJ/IZ
 Sample : H1256-18
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 9024

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-BF020116.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	4.167	179	182	187	rBV	298284	458306	3.23%	0.194%
2	4.876	240	244	246	rBV	3023339	4593482	32.41%	1.940%
3	5.310	279	282	284	rBV	3713043	4013206	28.32%	1.695%
4	5.710	315	317	320	rBV	369045	344609	2.43%	0.146%
5	6.362	371	374	378	rBV2	2959770	4835377	34.12%	2.042%
6	6.499	380	386	389	rBV3	4726998	11068935	78.10%	4.674%
7	6.704	402	404	407	rBV	815584	1087877	7.68%	0.459%
8	6.887	415	420	422	rBV2	5920966	8612021	60.76%	3.637%
9	6.990	426	429	430	rBV	2491350	3349619	23.63%	1.414%
10	7.150	438	443	446	rBV2	6927576	10297622	72.66%	4.348%
11	7.230	447	450	452	rVV	3511864	4029179	28.43%	1.701%
12	7.310	452	457	460	rVB2	4098765	8007355	56.50%	3.381%
13	7.550	475	478	482	rBV	3469243	4743038	33.47%	2.003%
14	7.630	482	485	487	rBV	2576817	3854861	27.20%	1.628%
15	7.767	493	497	499	rBV	3403127	3940683	27.80%	1.664%
16	7.870	503	506	508	rVV	3760221	3717068	26.23%	1.570%
17	7.939	510	512	515	rVV	1556271	2173031	15.33%	0.918%
18	8.030	515	520	522	rVB2	7129991	10039291	70.83%	4.239%
19	8.145	527	530	532	rBV	3409319	3311800	23.37%	1.398%
20	8.567	564	567	569	rBV	5764261	5957909	42.04%	2.516%
21	8.990	600	604	606	rBV	1565582	2393842	16.89%	1.011%
22	9.036	606	608	609	rVV	4445620	4134815	29.17%	1.746%
23	9.082	609	612	614	rVB	4660688	4752206	33.53%	2.007%
24	9.208	619	623	625	rVB	5608125	7090733	50.03%	2.994%
25	9.470	645	646	648	rVB	307983	349360	2.46%	0.148%
26	9.688	662	665	668	rBV2	122830	204748	1.44%	0.086%
27	9.745	668	670	672	rVV2	1411268	1607592	11.34%	0.679%
28	9.950	685	688	691	rBV2	4522924	6377288	45.00%	2.693%
29	10.190	705	709	711	rBV	6085230	6918578	48.82%	2.922%
30	10.305	715	719	722	rVB	11399308	14173056	100.00%	5.985%
31	10.431	725	730	732	rBV2	175577	253383	1.79%	0.107%
32	10.533	737	739	742	rVB2	2179811	2878204	20.31%	1.215%
33	10.659	749	750	755	rBV	158400	242841	1.71%	0.103%
34	10.785	757	761	763	rBV	5441139	6195641	43.71%	2.616%

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Sampling : 1 Min Area: 3 % of largest Peak
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Stop Thrs : 0 Peak Location: TOP

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35	10.853	763	767	769	rBV	2342241	3398266	23.98%	1.435%
36	11.048	781	784	786	rBV	4197455	4587167	32.37%	1.937%
37	11.105	788	789	795	rVB2	105890	221766	1.56%	0.094%
38	11.231	797	800	801	rVV	1762553	1954048	13.79%	0.825%
39	11.253	801	802	804	rVV	3482703	3443650	24.30%	1.454%
40	11.311	804	807	809	rVB	5611568	5505745	38.85%	2.325%
41	11.471	818	821	825	rBV	236243	363333	2.56%	0.153%
42	12.294	891	893	898	rBV	407580	694153	4.90%	0.293%
43	12.442	903	906	909	rBV	5468990	7601901	53.64%	3.210%
44	12.671	923	926	929	rBV	5570055	6842267	48.28%	2.889%
45	12.819	936	939	941	rBV	4490736	4686609	33.07%	1.979%
46	13.299	978	981	984	rVB	5951327	7256689	51.20%	3.064%
47	13.722	1016	1018	1027	rVB	321875	521000	3.68%	0.220%
48	13.859	1027	1030	1031	rBV2	5185107	6111098	43.12%	2.581%
49	13.894	1031	1033	1035	rVB	1959852	2560302	18.06%	1.081%
50	14.362	1070	1074	1081	rBV3	65675	239075	1.69%	0.101%
51	14.477	1081	1084	1086	rBV	6571018	7998110	56.43%	3.377%
52	14.705	1100	1104	1107	rVB	139382	150455	1.06%	0.064%
53	14.945	1122	1125	1127	rBV	138799	166684	1.18%	0.070%
54	15.185	1143	1146	1149	rBV	1590331	1971414	13.91%	0.832%
55	15.242	1149	1151	1153	rVB	866576	940195	6.63%	0.397%
56	15.677	1186	1189	1192	rBV	112048	179504	1.27%	0.076%
57	16.580	1261	1268	1271	rBV2	2644702	6694592	47.23%	2.827%
58	16.980	1295	1303	1306	rBV	2328654	6198880	43.74%	2.618%
59	17.083	1309	1312	1318	rVB7	51308	181972	1.28%	0.077%
60	17.974	1385	1390	1395	rVB3	40833	143829	1.01%	0.061%
61	18.866	1462	1468	1474	rVB5	59790	192068	1.36%	0.081%

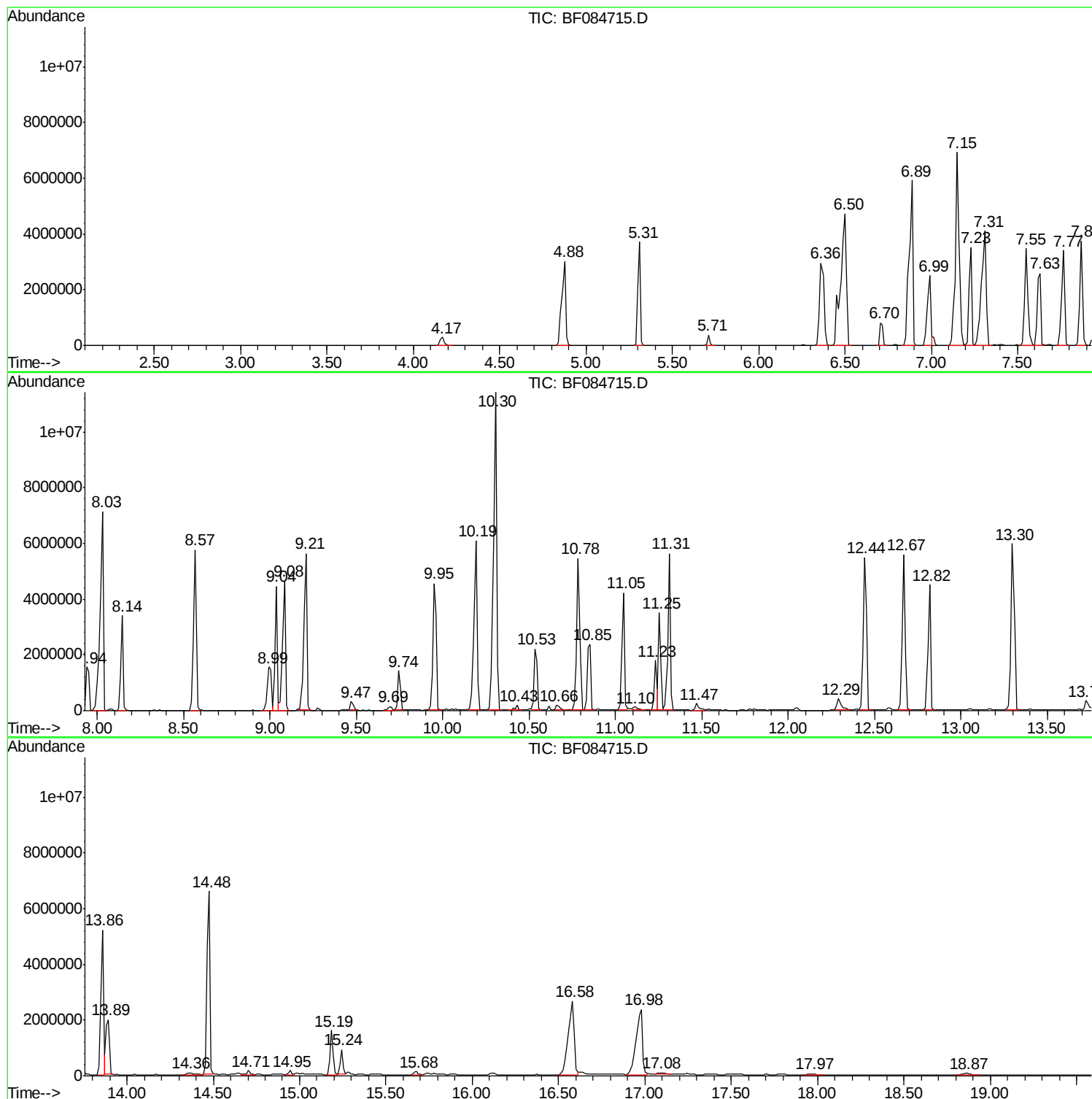
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF020116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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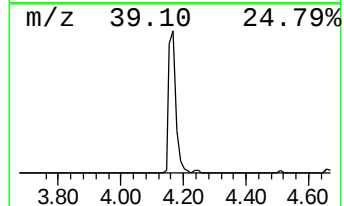
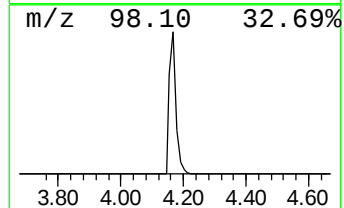
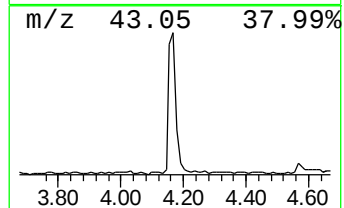
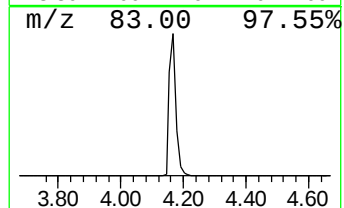
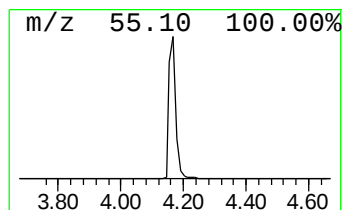
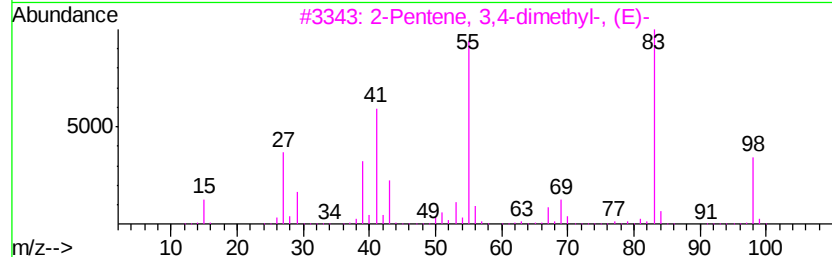
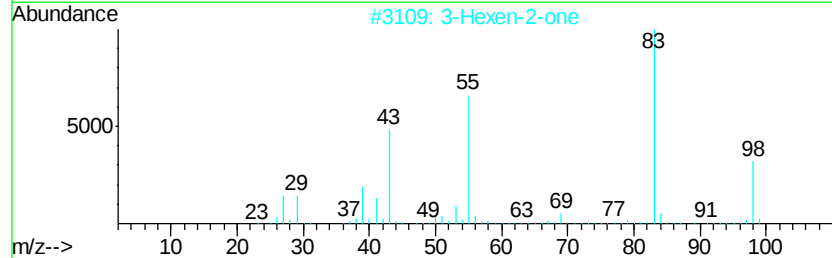
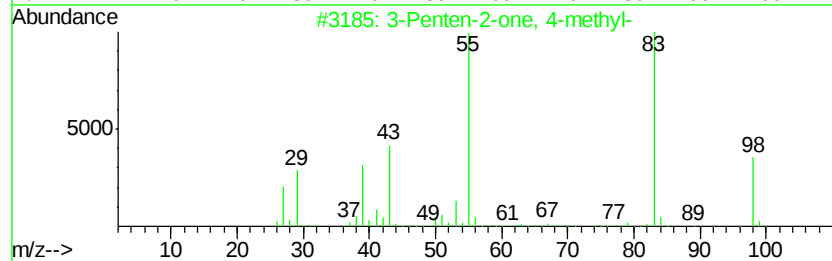
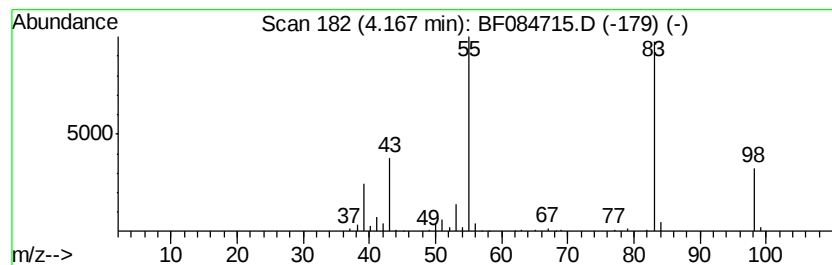
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF020116.M
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.17	8.43 ng	458306	1,4-Dichlorobenzene-d4	6.72

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Penten-2-one, 4-methyl-	98	C6H10O	000141-79-7	94
2			3-Hexen-2-one	98	C6H10O	000763-93-9	90
3			2-Pentene, 3,4-dimethyl-, (E)-	98	C7H14	004914-92-5	90
4			2-Pentene, 3,4-dimethyl-, (Z)-	98	C7H14	004914-91-4	86
5			2-Pentene, 4,4-dimethyl-, (E)-	98	C7H14	000690-08-4	86



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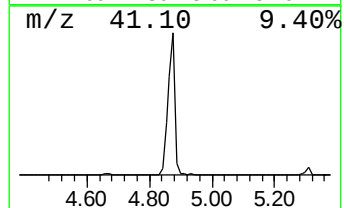
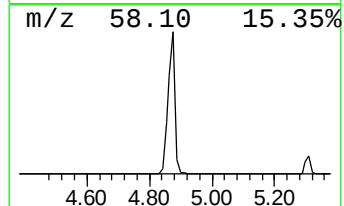
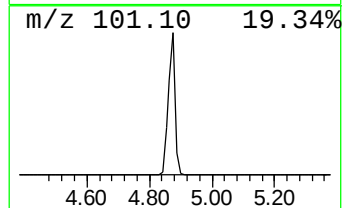
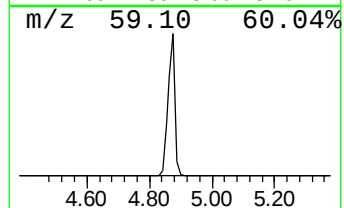
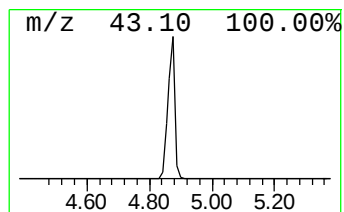
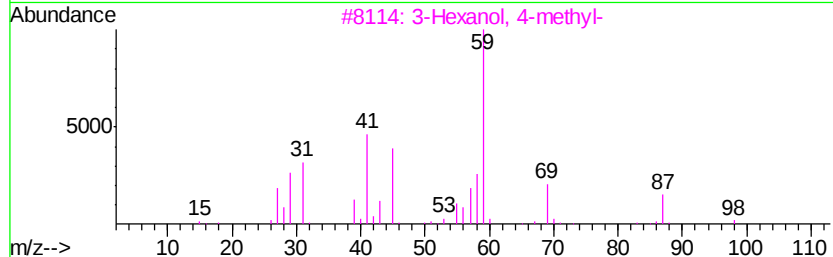
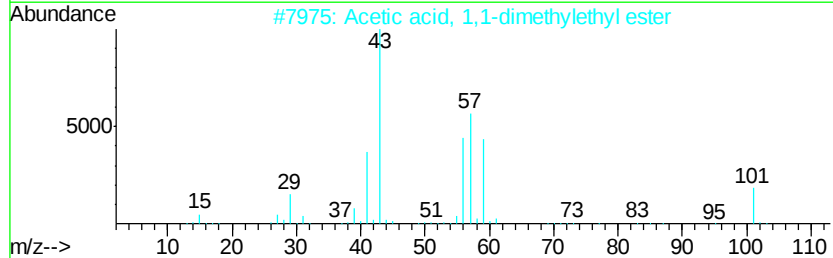
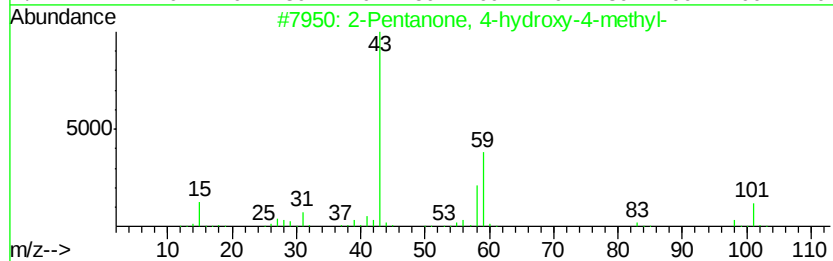
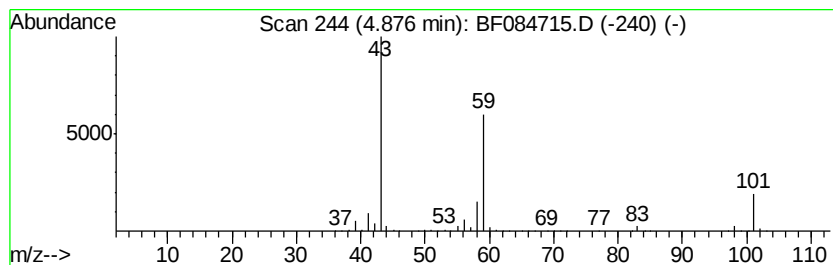
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.88	84.45 ng	4593480	1,4-Dichlorobenzene-d4	6.72

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2		Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
3		3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	28
4		Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	25
5		Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9



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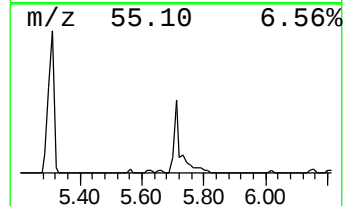
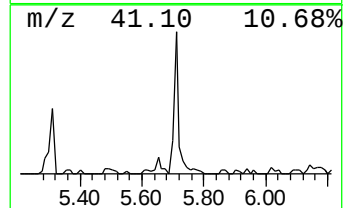
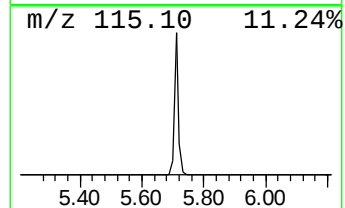
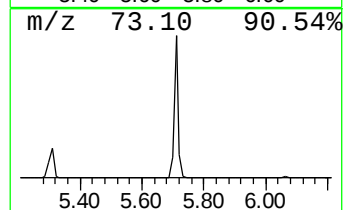
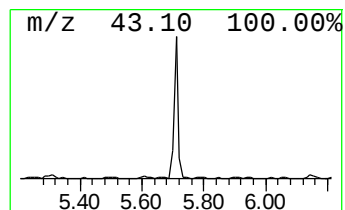
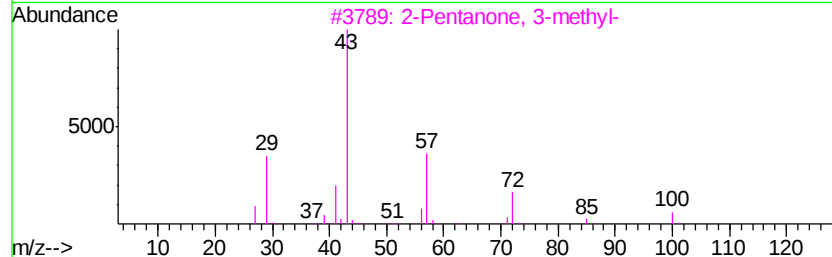
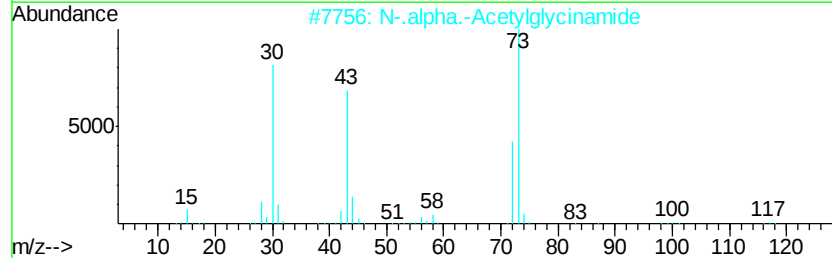
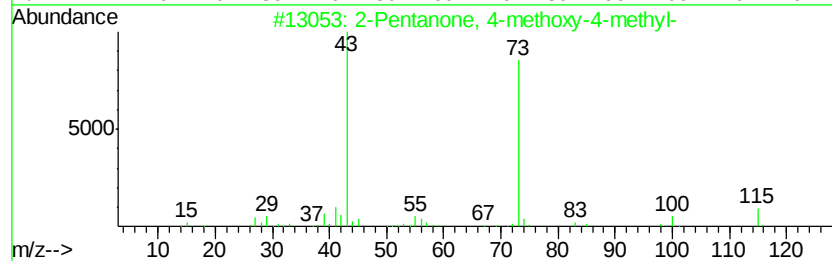
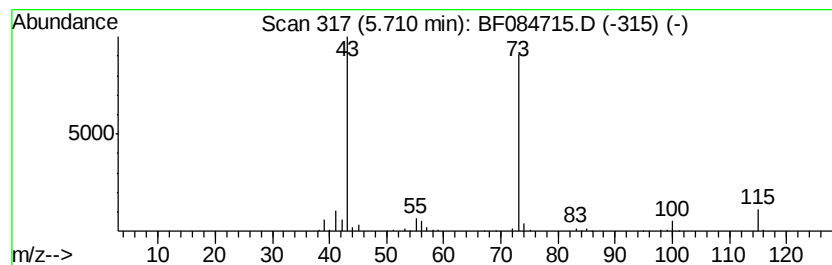
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 3 2-Pentanone, 4-methoxy-4-me... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.71	6.34 ng	344609	1,4-Dichlorobenzene-d4	6.72

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-methoxy-4-methyl-	130	C7H14O2	000107-70-0	83
2		N-.alpha.-Acetylglucineamide	116	C4H8N2O2	002620-63-5	38
3		2-Pentanone, 3-methyl-	100	C6H12O	000565-61-7	16
4		N-Formylmorpholine	115	C5H9NO2	004394-85-8	10
5		2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	10



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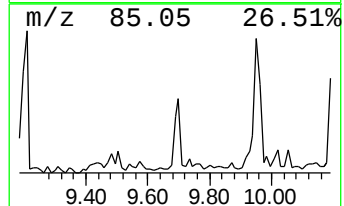
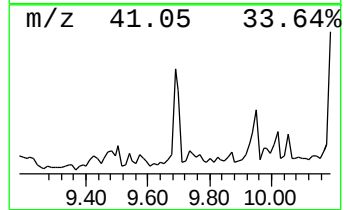
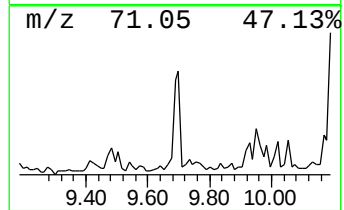
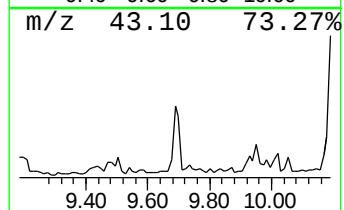
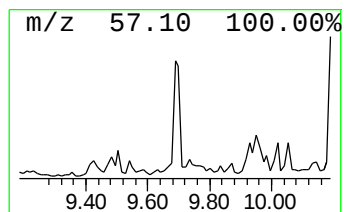
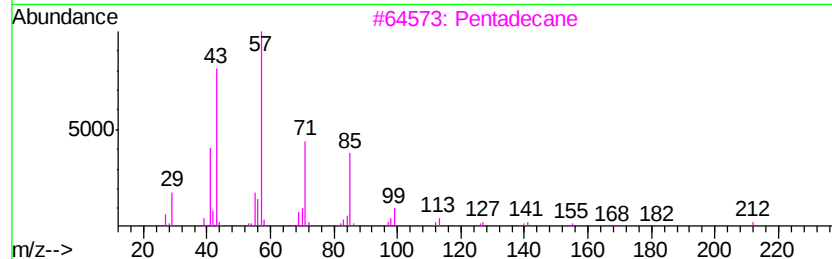
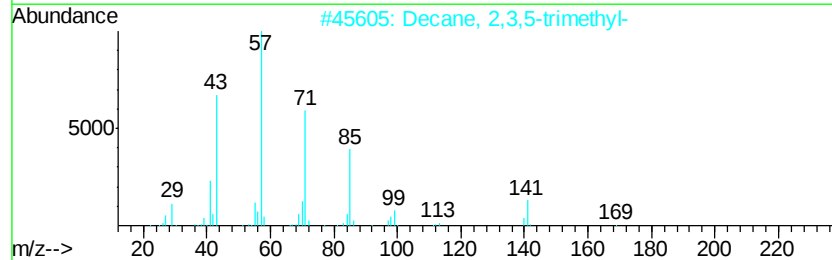
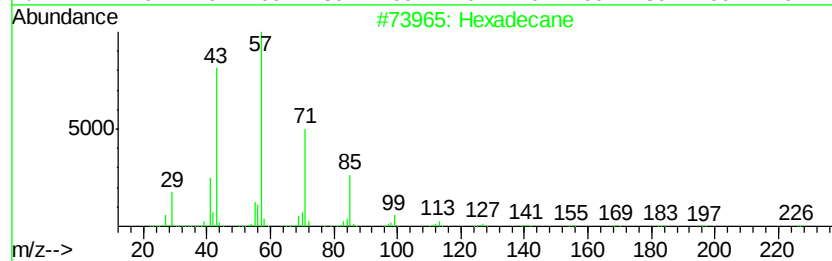
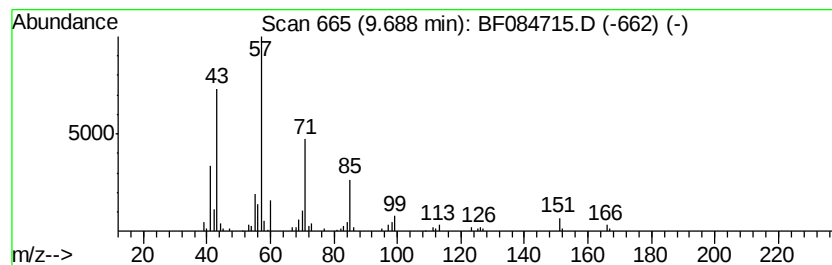
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 4 Hexadecane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.69	2.55 ng	204748	Acenaphthene-d10	9.74

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexadecane	226	C16H34	000544-76-3	74
2			Decane, 2,3,5-trimethyl-	184	C13H28	062238-11-3	72
3			Pentadecane	212	C15H32	000629-62-9	64
4			Tetradecane	198	C14H30	000629-59-4	64
5			Eicosane	282	C20H42	000112-95-8	58



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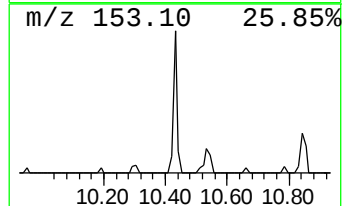
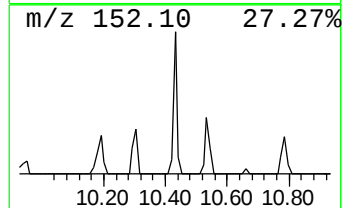
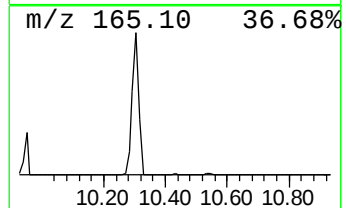
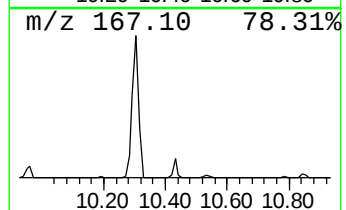
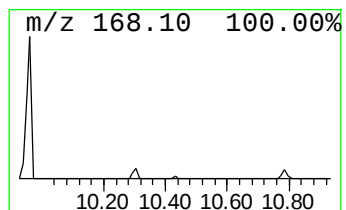
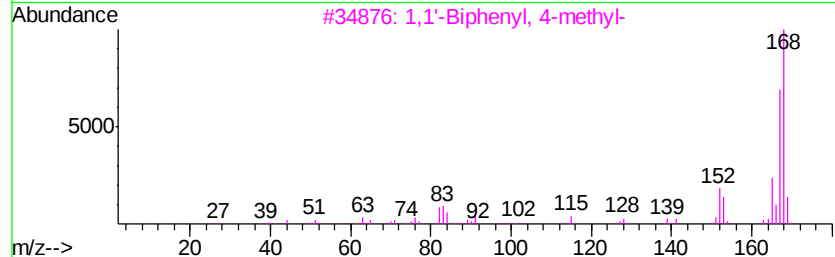
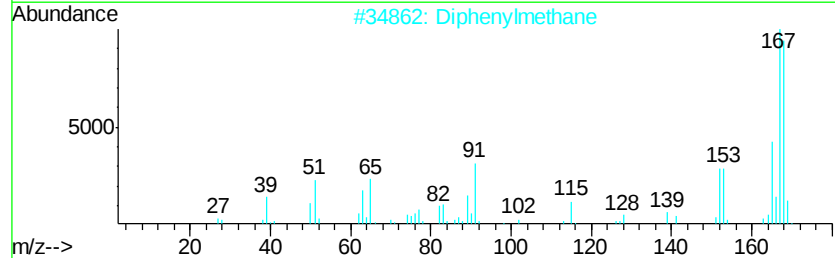
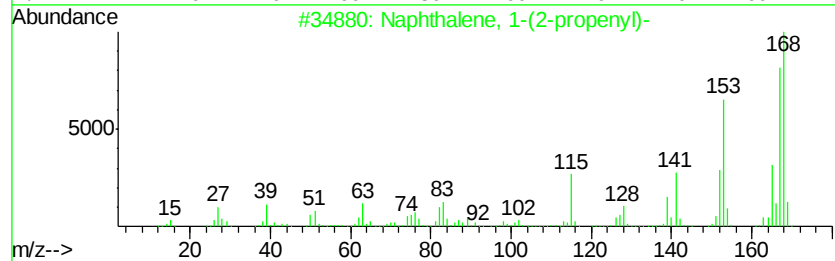
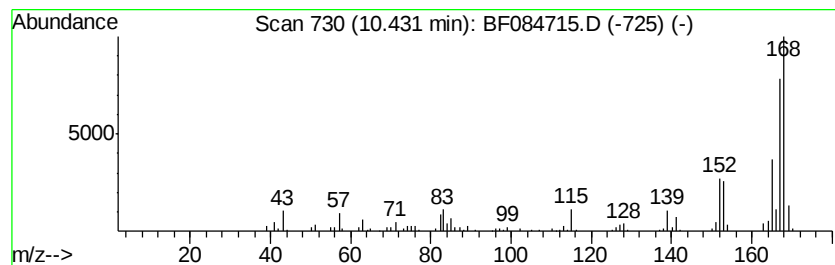
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF020116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Naphthalene, 1-(2-propenyl)- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.43	3.15 ng	253383	Acenaphthene-d10	9.74
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 1-(2-propenyl)-	168 C13H12	002489-86-3 93
2		Diphenylmethane	168 C13H12	000101-81-5 93
3		1,1'-Biphenyl, 4-methyl-	168 C13H12	000644-08-6 93
4		1,1'-Biphenyl, 2-methyl-	168 C13H12	000643-58-3 91
5		1,1'-Biphenyl, 3-methyl-	168 C13H12	000643-93-6 89



Data Path : Z:\HPCHEM1\BNA F\DATA\BF020116\
Data File : BF084715.D
Acq On : 1 Feb 2016 18:00
Operator : SJ/IZ
Sample : H1256-18
Misc :
ALS Vial : 16 Sample Multiplier: 1

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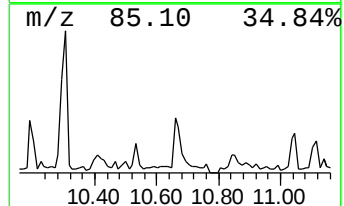
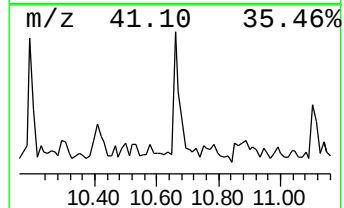
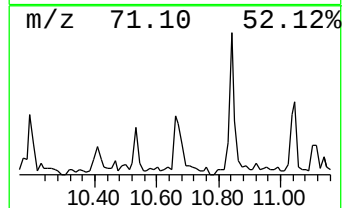
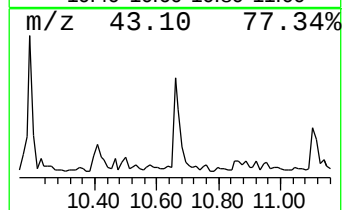
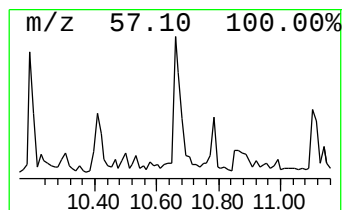
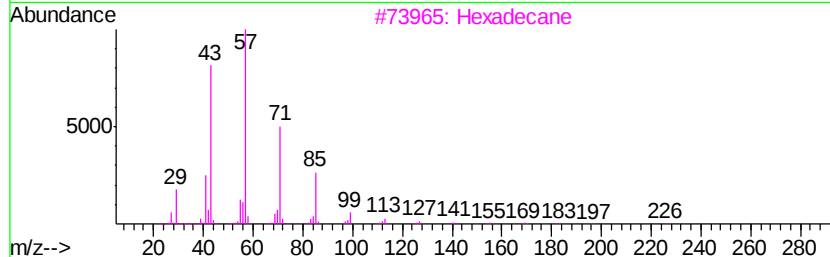
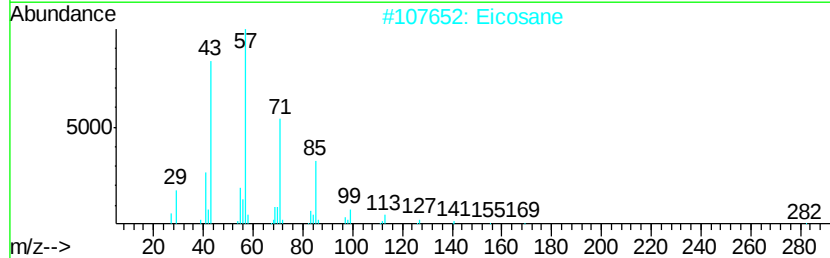
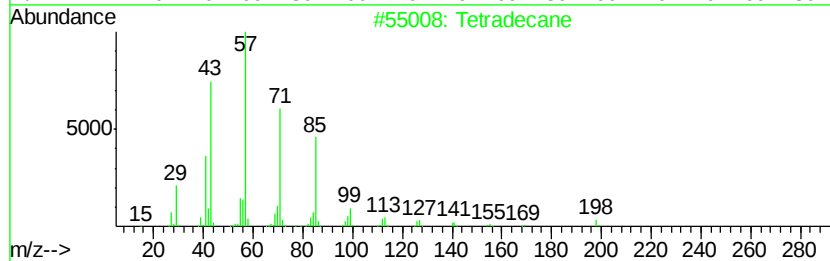
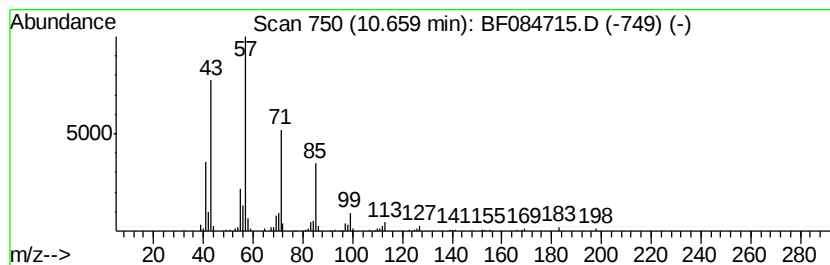
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Tetradecane Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.66	2.49 ng	242841	Phenanthrene-d10	11.23

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tetradecane	198	C14H30	000629-59-4	94
2		Eicosane	282	C20H42	000112-95-8	91
3		Hexadecane	226	C16H34	000544-76-3	91
4		Heptadecane	240	C17H36	000629-78-7	90
5		Pentadecane	212	C15H32	000629-62-9	86



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF020116\
Data File : BF084715.D
Acq On : 1 Feb 2016 18:00
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Sample : H1256-18
Misc :
ALS Vial : 16 Sample Multiplier: 1

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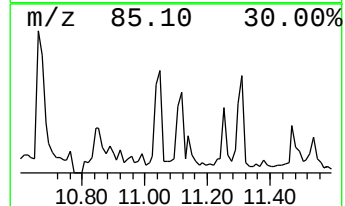
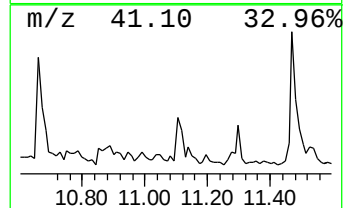
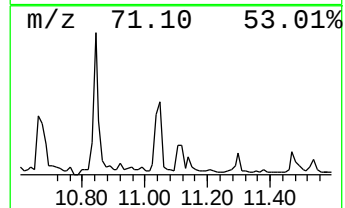
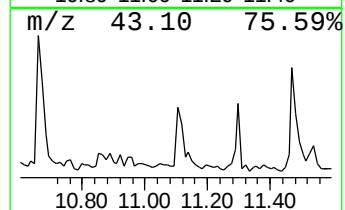
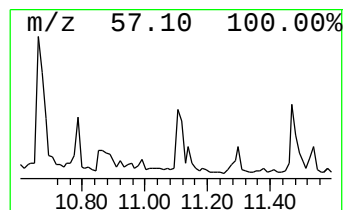
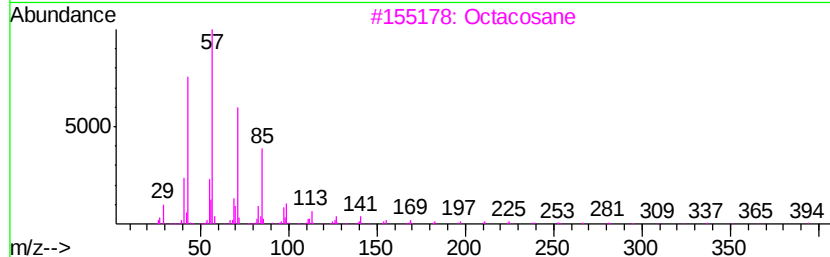
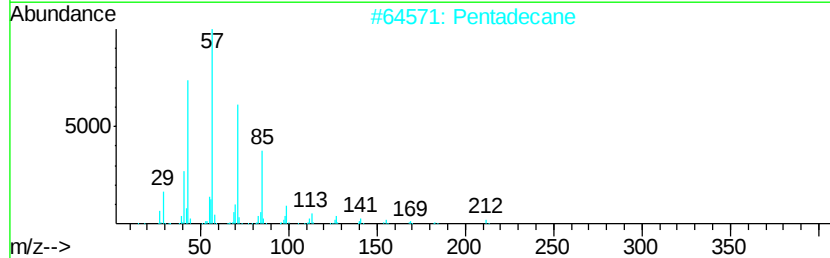
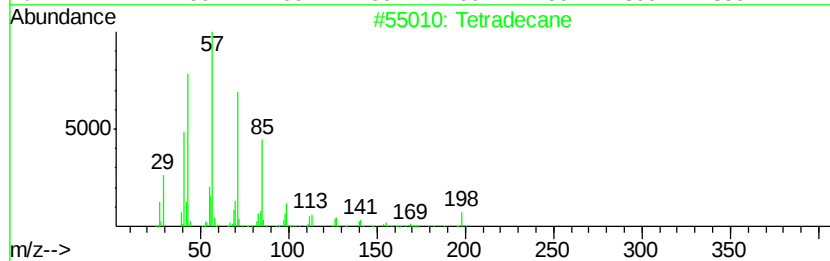
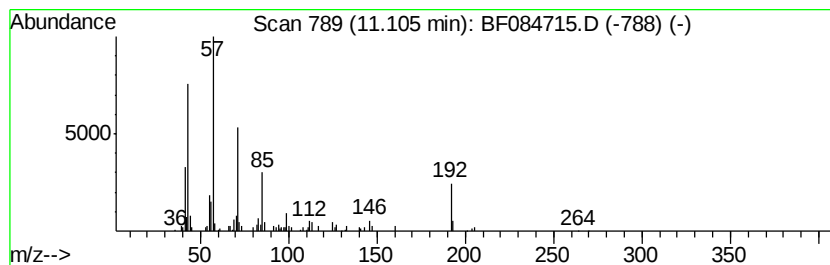
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Pentadecane Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.10	2.27 ng	221766	Phenanthrene-d10	11.23

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tetradecane	198	C14H30	000629-59-4	59
2		Pentadecane	212	C15H32	000629-62-9	59
3		Octacosane	394	C28H58	000630-02-4	59
4		Hexadecane	226	C16H34	000544-76-3	58
5		2-Bromo dodecane	248	C12H25Br	013187-99-0	53



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF020116\
Data File : BF084715.D
Acq On : 1 Feb 2016 18:00
Operator : SJ/IZ
Sample : H1256-18
Misc :
ALS Vial : 16 Sample Multiplier: 1

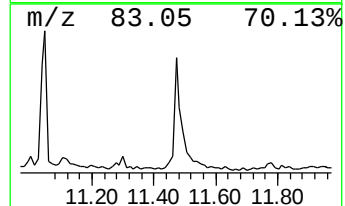
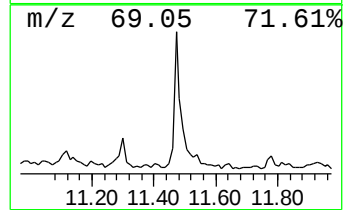
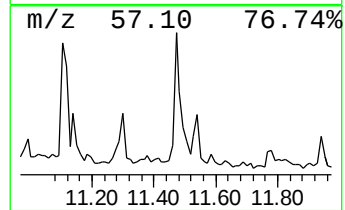
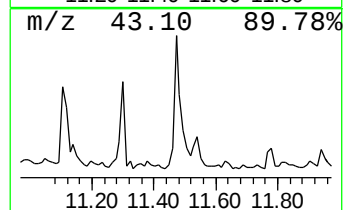
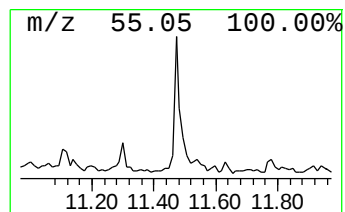
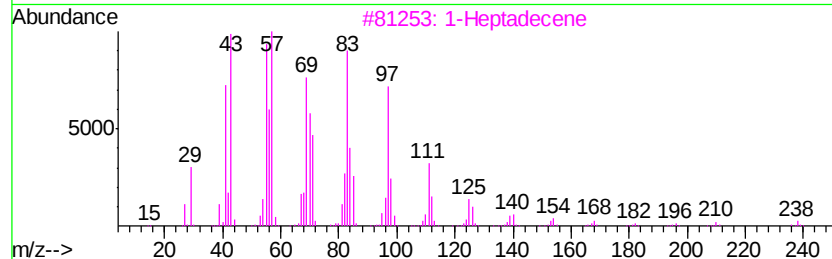
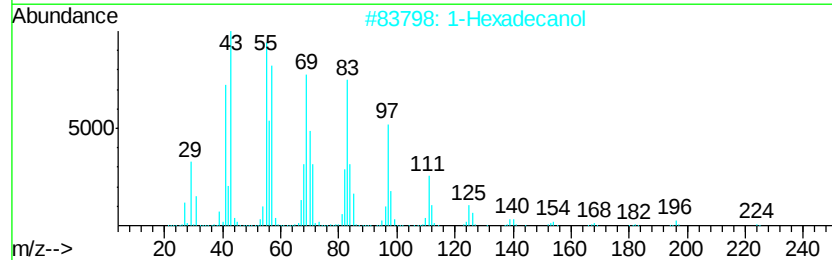
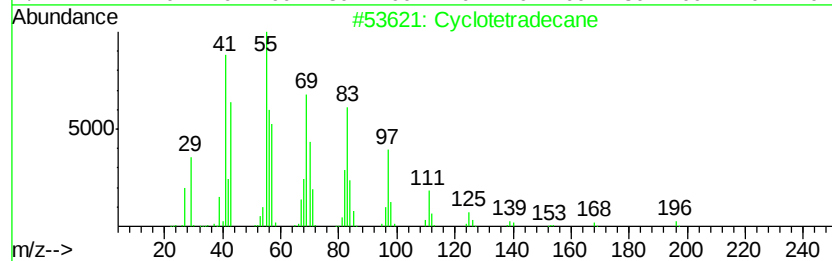
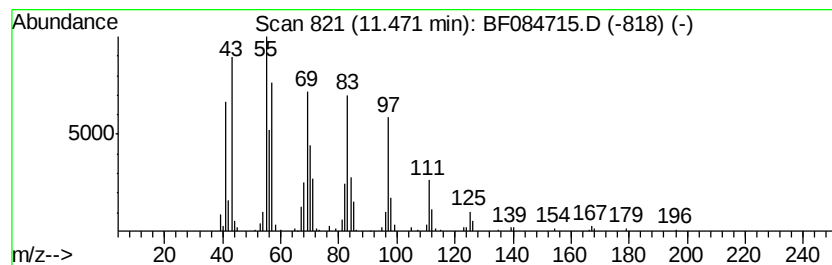
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Cyclotetradecane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD		R.T.
11.47	3.72 ng	363333	Phenanthrene-d10		11.23
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclotetradecane	196	C14H28	000295-17-0	95
2	1-Hexadecanol	242	C16H34O	036653-82-4	94
3	1-Heptadecene	238	C17H34	006765-39-5	91
4	1-Hexadecene	224	C16H32	000629-73-2	91
5	9-Octadecene, (E)-	252	C18H36	007206-25-9	91



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF020116\
Data File : BF084715.D
Acq On : 1 Feb 2016 18:00
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Sample : H1256-18
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ClientSampleId :
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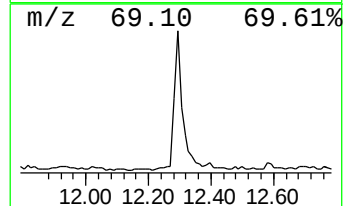
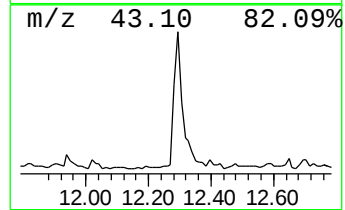
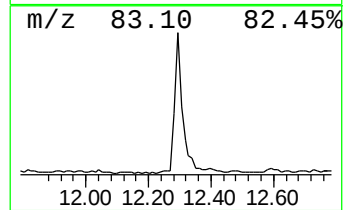
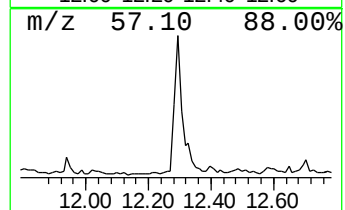
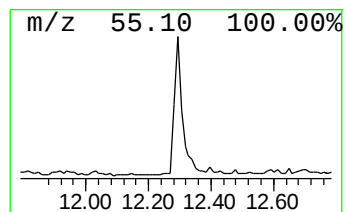
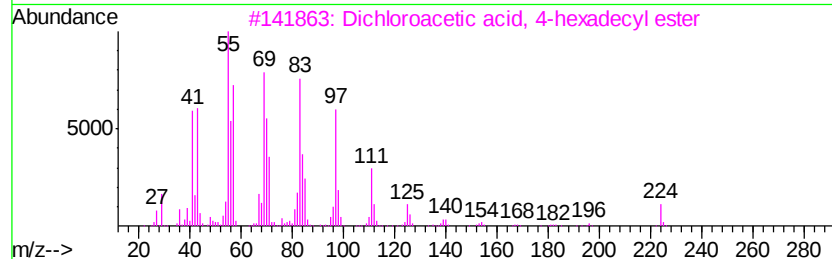
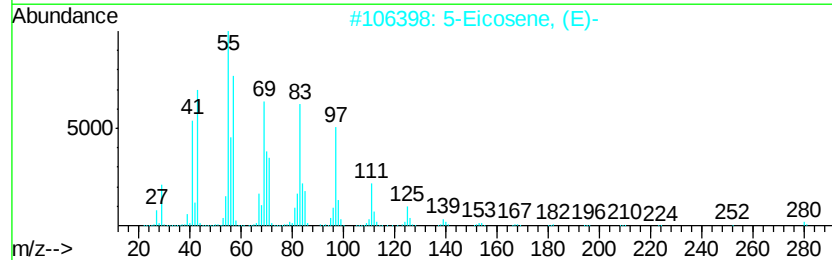
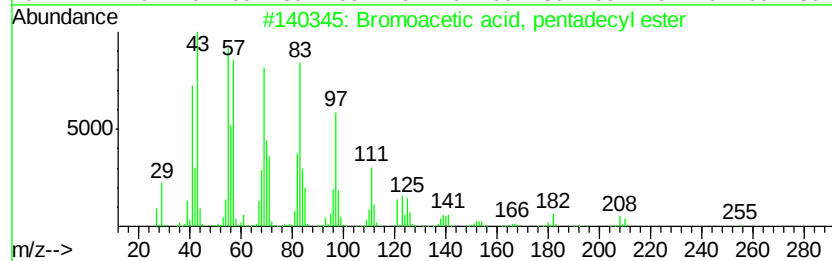
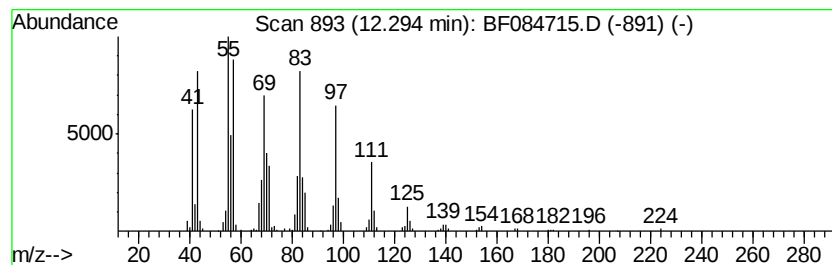
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Bromoacetic acid, pentadecyl... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.29	7.10 ng	694153	Phenanthrene-d10	11.23

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Bromoacetic acid, pentadecyl ester	348	C17H33BrO2	131143-01-6	91
2		5-Eicosene, (E)-	280	C20H40	074685-30-6	91
3		Dichloroacetic acid, 4-hexadecyl...	352	C18H34Cl2O2	1000282-98-1	91
4		Cyclohexadecane	224	C16H32	000295-65-8	91
5		1-Nonadecene	266	C19H38	018435-45-5	91



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF020116\
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Acq On : 1 Feb 2016 18:00
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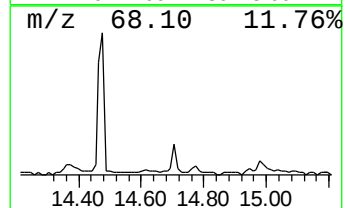
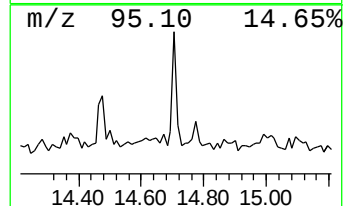
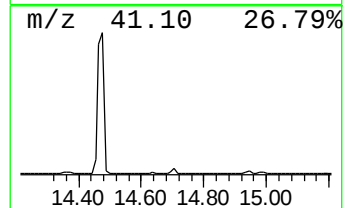
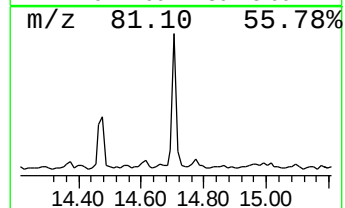
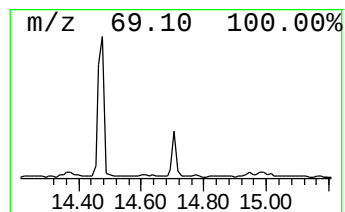
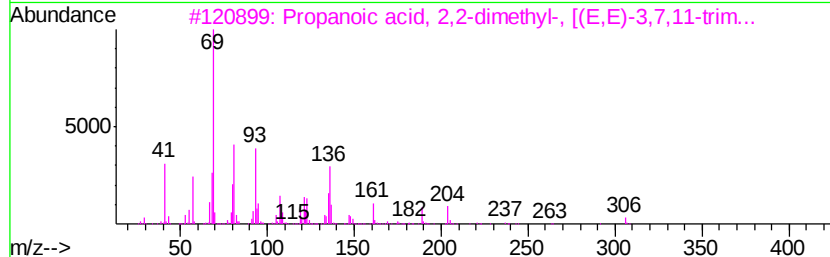
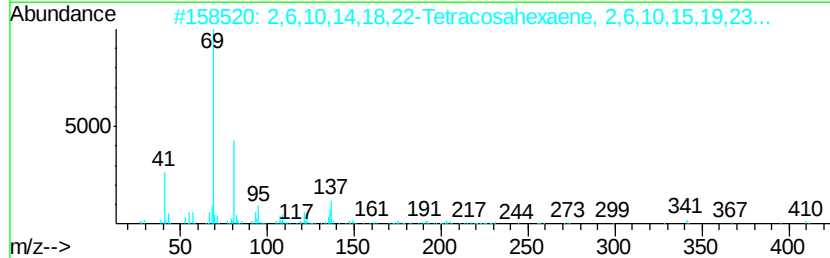
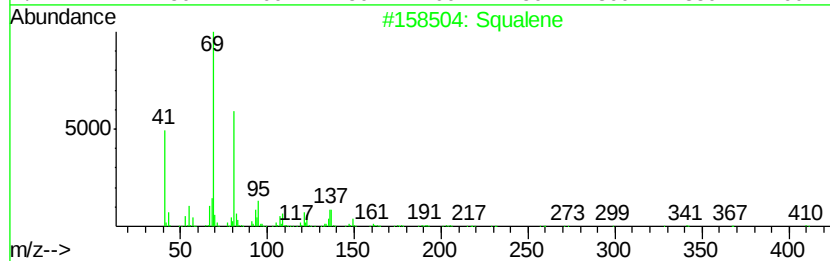
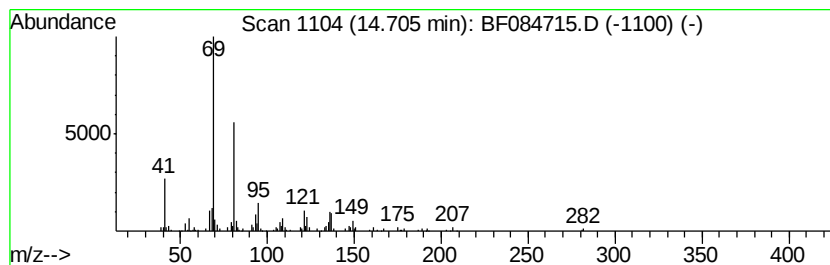
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Squalene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.71	3.20 ng	150455	Perylene-d12	15.24

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Squalene	410	C30H50	007683-64-9	91
2		2,6,10,14,18,22-Tetracosahexaene...	410	C30H50	000111-02-4	86
3		Propanoic acid, 2,2-dimethyl-, [...	306	C20H34O2	1000164-38-8	64
4		7,11-Dimethyldodeca-2,6,10-trien...	208	C14H24	125529-10-4	64
5		1,5,9-Decatriene, 2,3,5,8-tetram...	192	C14H24	230646-72-7	59



Data Path : Z:\HPCHEM1\BNA F\DATA\BF020116\
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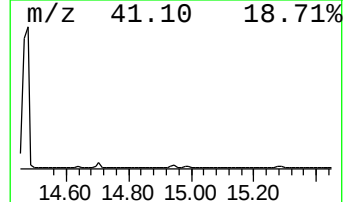
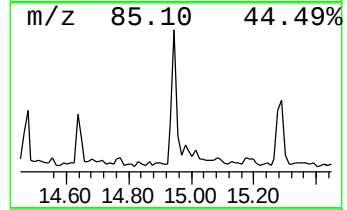
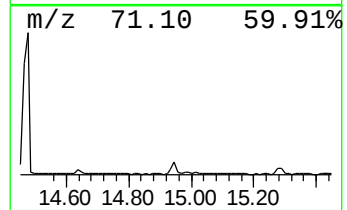
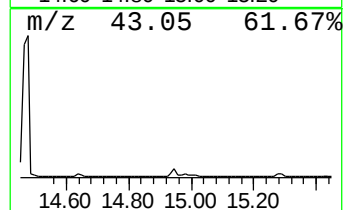
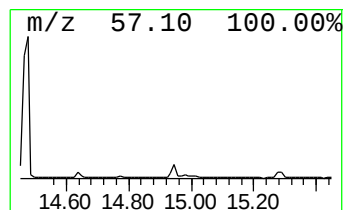
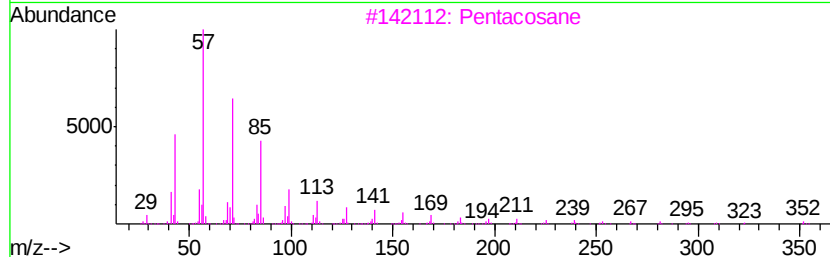
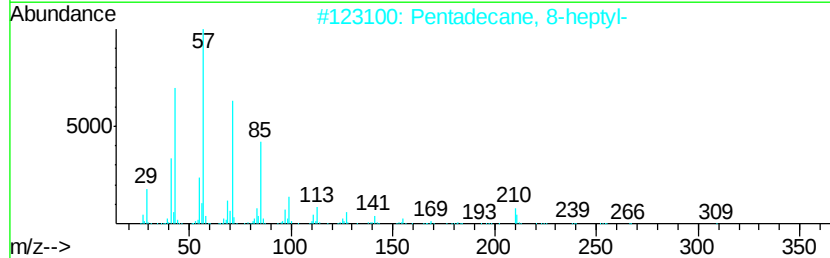
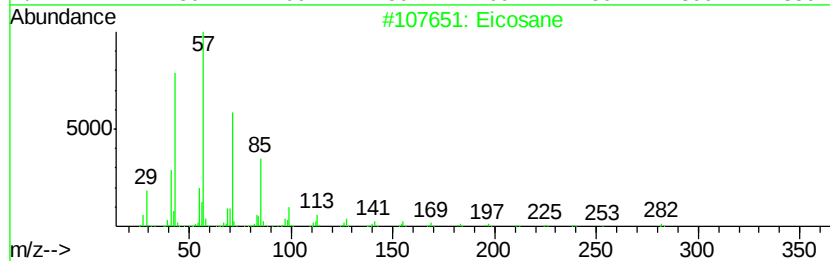
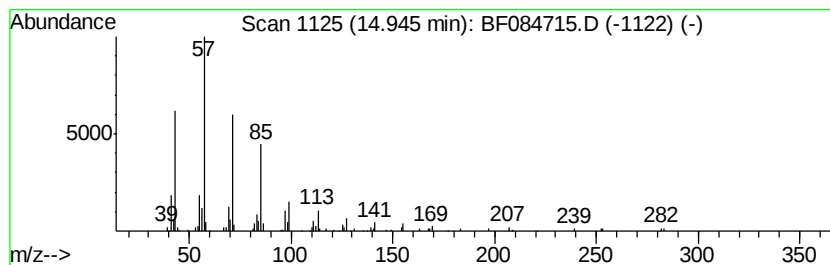
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Eicosane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.95	3.55 ng	166684	Perylene-d12	15.24

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	96
2		Pentadecane, 8-heptyl-	310	C22H46	071005-15-7	91
3		Pentacosane	352	C25H52	000629-99-2	91
4		Nonacosane	408	C29H60	000630-03-5	91
5		Octacosane	394	C28H58	000630-02-4	91



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ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
9024

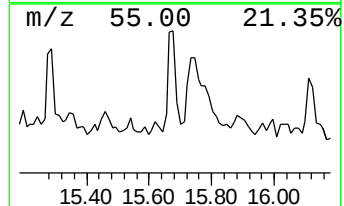
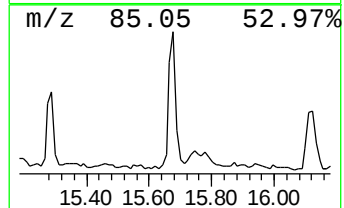
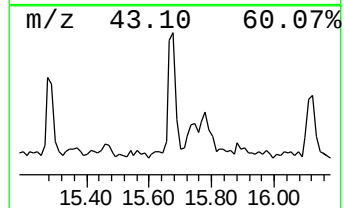
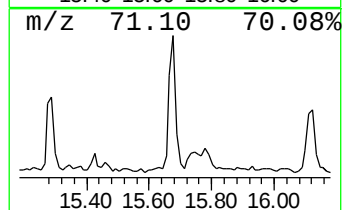
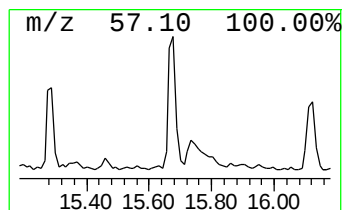
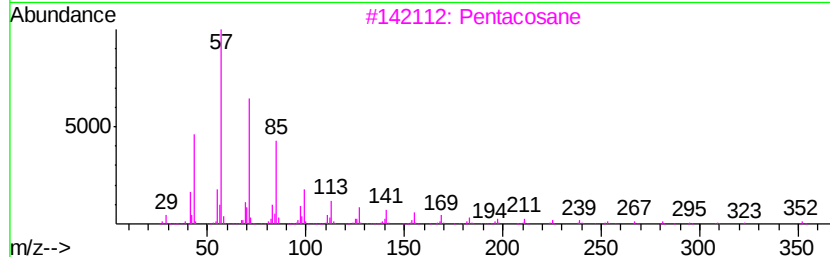
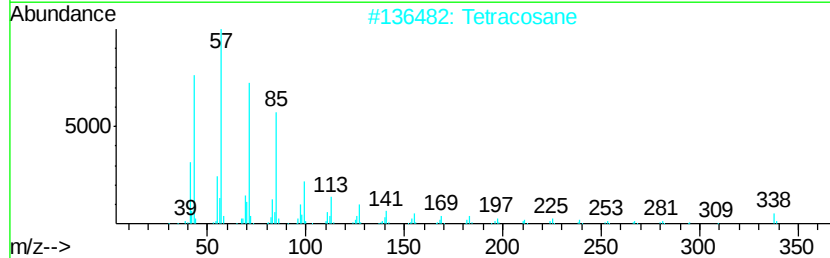
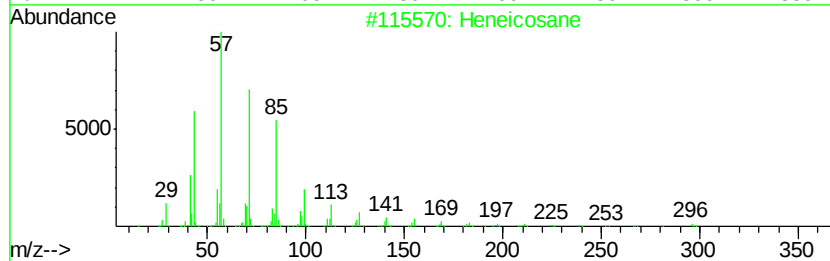
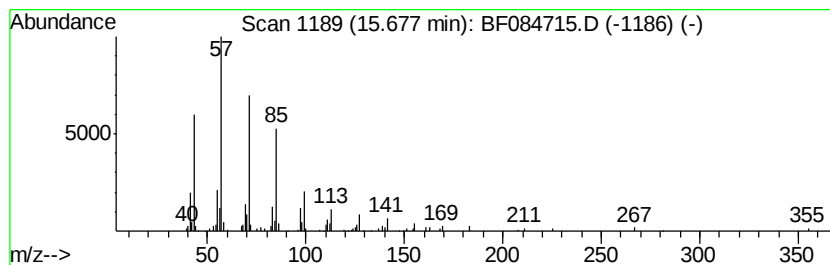
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Heneicosane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.68	3.82 ng	179504	Perylene-d12	15.24

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	91
2		Tetracosane	338	C24H50	000646-31-1	91
3		Pentacosane	352	C25H52	000629-99-2	91
4		Pentadecane, 8-heptyl-	310	C22H46	071005-15-7	90
5		triacontane	422	C30H62	000638-68-6	87



Data Path : Z:\HPCHEM1\BNA F\DATA\BF020116\
Data File : BF084715.D
Acq On : 1 Feb 2016 18:00
Operator : SJ/IZ
Sample : H1256-18
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
9024

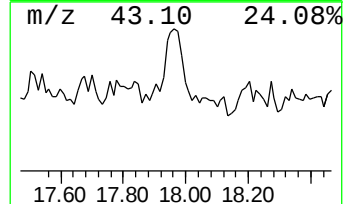
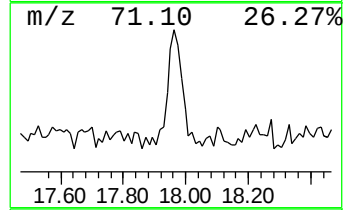
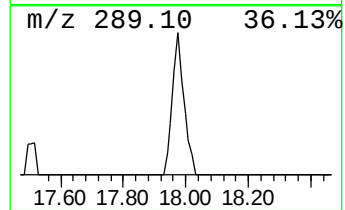
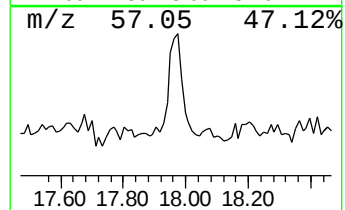
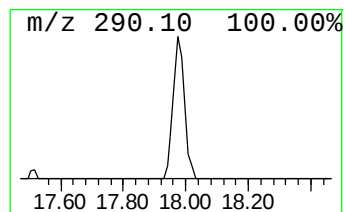
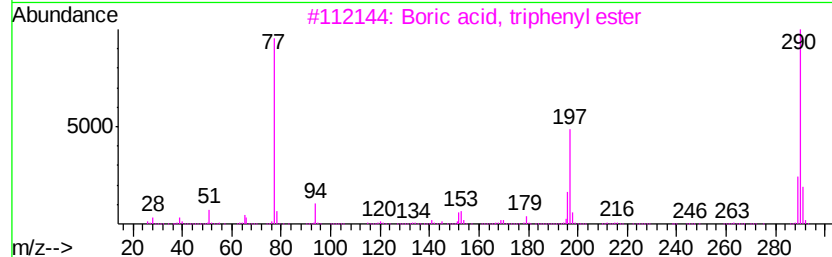
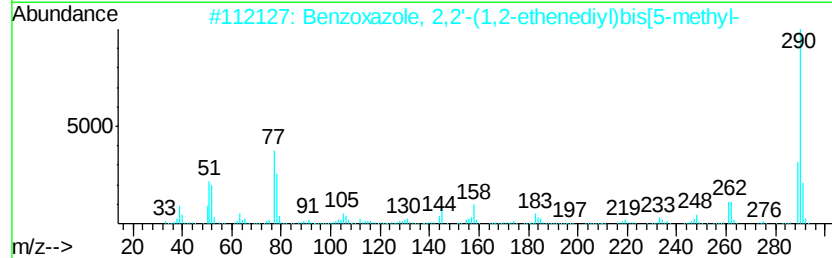
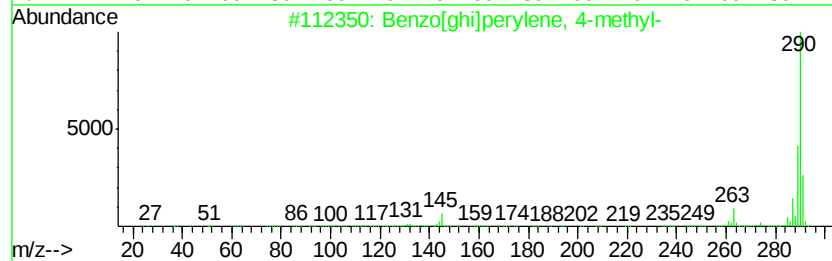
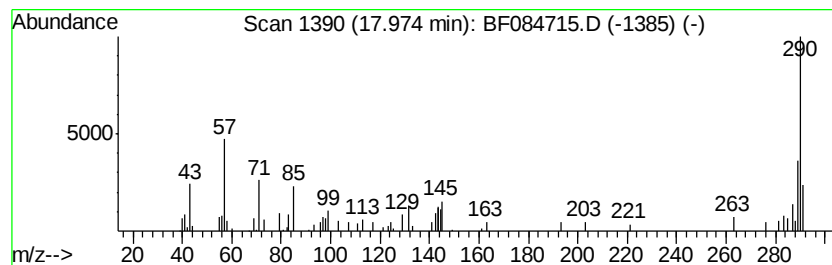
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Benzo[ghi]perylene, 4-methyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.97	3.06 ng	143829	Perylene-d12	15.24

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[ghi]perylene, 4-methyl-	290	C23H14	019224-38-5	58
2			Benzoxazole, 2,2'-(1,2-ethenediv...	290	C18H14N2O2	001041-00-5	40
3			Boric acid, triphenyl ester	290	C18H15B03	001095-03-0	38
4			5.alpha.-Androstan-17.beta.-ol, ...	290	C20H34O	005846-88-8	33
5			3-Trifluoromethyl-9-phenanthrene...	290	C16H9F3O2	035315-61-8	32



Data Path : Z:\HPCHEM1\BNA F\DATA\BF020116\
Data File : BF084715.D
Acq On : 1 Feb 2016 18:00
Operator : SJ/IZ
Sample : H1256-18
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
9024

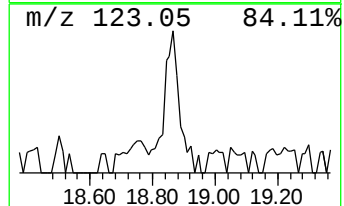
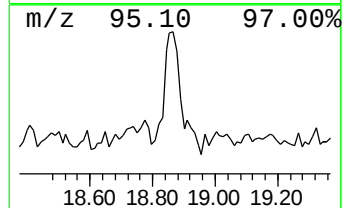
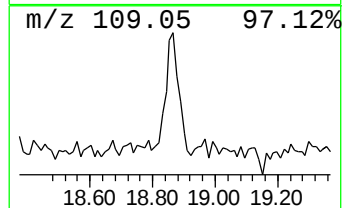
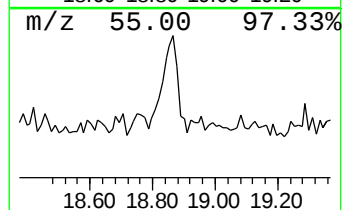
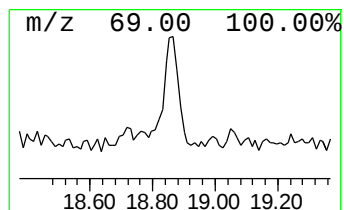
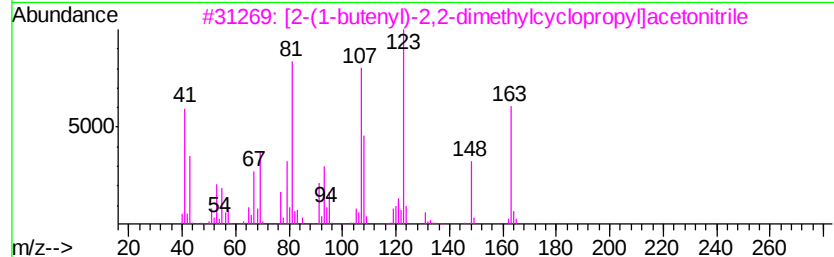
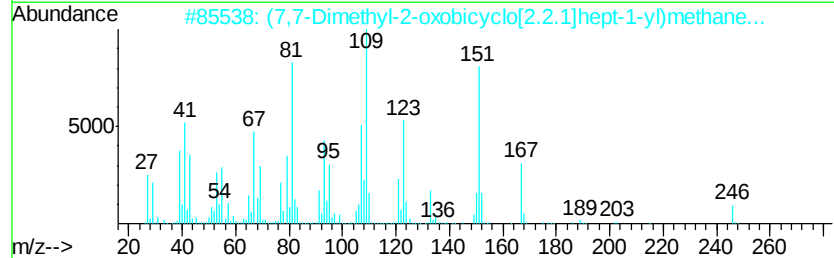
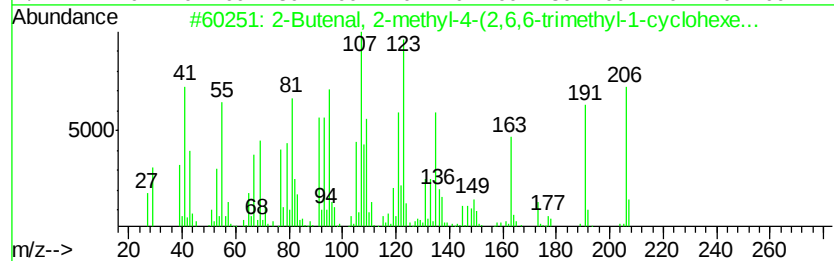
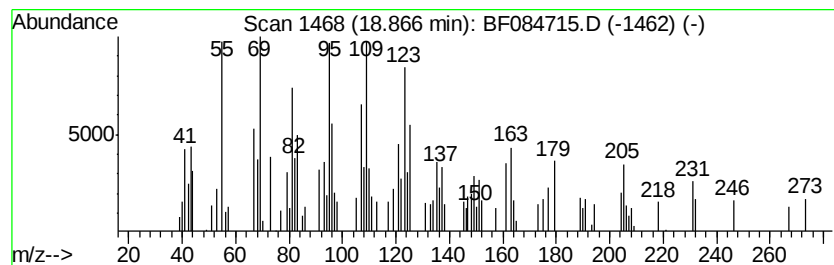
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 unknown18.87 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.87	4.09 ng	192068	Perylene-d12	15.24

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Butenal, 2-methyl-4-(2,6,6-tri...	206	C14H22O	003155-71-3	47
2		(7,7-Dimethyl-2-oxobicyclo[2.2.1...	246	C11H18O4S	1000197-55-6	46
3		[2-(1-butenyl)-2,2-dimethylcyclo...	163	C11H17N	1000111-15-4	42
4		3,4-Octadiene, 7-methyl-	124	C9H16	037050-05-8	25
5		But-3-enal, 2-methyl-4-(2,6,6-tr...	206	C14H22O	104900-59-6	20



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF020116\
 Data File : BF084715.D
 Acq On : 1 Feb 2016 18:00
 Operator : SJ/IJ
 Sample : H1256-18
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 9024

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF020116.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...	4.17	8.4	ng	458306	1	6.72	1087880	20.0
2-Pentanone, 4-hy...	4.88	84.5	ng	4593480	1	6.72	1087880	20.0
2-Pentanone, 4-me...	5.71	6.3	ng	344609	1	6.72	1087880	20.0
Hexadecane	9.69	2.5	ng	204748	3	9.74	1607590	20.0
Naphthalene, 1-(2...	10.43	3.1	ng	253383	3	9.74	1607590	20.0
Tetradecane	10.66	2.5	ng	242841	4	11.23	1954050	20.0
Pentadecane	11.10	2.3	ng	221766	4	11.23	1954050	20.0
Cyclotetradecane	11.47	3.7	ng	363333	4	11.23	1954050	20.0
Bromoacetic acid,...	12.29	7.1	ng	694153	4	11.23	1954050	20.0
Squalene	14.71	3.2	ng	150455	6	15.24	940195	20.0
Eicosane	14.95	3.5	ng	166684	6	15.24	940195	20.0
Heneicosane	15.68	3.8	ng	179504	6	15.24	940195	20.0
Benzo[ghi]perylene...	17.97	3.1	ng	143829	6	15.24	940195	20.0
unknown18.87	18.87	4.1	ng	192068	6	15.24	940195	20.0