

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF120815\
 Data File : BF083631.D
 Acq On : 9 Dec 2015 9:39
 Operator : UM/IZ
 Sample : G4647-19
 Misc :
 ALS Vial : 39 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 BRIDGE14-10(0-2)

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-BF120815.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	3.538	160	162	176	rBV	365129	753526	23.79%	2.512%
2	3.984	199	201	218	rBV	2167858	3002743	94.82%	10.012%
3	5.276	311	314	323	rBV	1850343	3166840	100.00%	10.559%
4	5.401	323	325	337	rVB	2422922	3156520	99.67%	10.524%
5	5.699	349	351	359	rBV	643239	808713	25.54%	2.696%
6	5.916	368	370	381	rVB	1723141	2159328	68.19%	7.200%
7	6.327	404	406	419	rVB	36123	83245	2.63%	0.278%
8	6.533	421	424	438	rBV	1162072	1976283	62.41%	6.589%
9	7.607	515	518	530	rBV	602838	968888	30.59%	3.230%
10	9.150	651	653	657	rBV	19869	39686	1.25%	0.132%
11	9.345	667	670	673	rBV	1875068	2434891	76.89%	8.118%
12	9.562	684	689	692	rVB2	22198	42013	1.33%	0.140%
13	10.030	727	730	733	rBV	376884	466042	14.72%	1.554%
14	10.213	744	746	752	rVB2	21732	36024	1.14%	0.120%
15	10.385	758	761	766	rBV2	611706	1034300	32.66%	3.449%
16	10.796	792	797	799	rBV2	14384	33562	1.06%	0.112%
17	10.933	807	809	812	rVB	24356	40446	1.28%	0.135%
18	11.231	832	835	838	rVV	113975	144458	4.56%	0.482%
19	11.596	862	867	869	rBV3	32932	81136	2.56%	0.271%
20	11.688	872	875	881	rVV	857692	1455392	45.96%	4.852%
21	11.779	881	883	887	rVB2	22180	38474	1.21%	0.128%
22	12.008	900	903	907	rBV	95491	198004	6.25%	0.660%
23	12.076	907	909	911	rVV2	26960	38963	1.23%	0.130%
24	12.316	923	930	932	rBV5	19116	53335	1.68%	0.178%
25	12.419	937	939	942	rBV	25009	42799	1.35%	0.143%
26	12.739	965	967	969	rBV	49237	61545	1.94%	0.205%
27	12.785	969	971	978	rVB	609808	849923	26.84%	2.834%
28	13.425	1025	1027	1032	rBV2	38116	68807	2.17%	0.229%
29	13.711	1050	1052	1055	rBV3	17047	34106	1.08%	0.114%
30	13.768	1055	1057	1060	rVV4	27327	47813	1.51%	0.159%
31	13.837	1060	1063	1071	rVV	134793	228160	7.20%	0.761%
32	14.214	1092	1096	1098	rBV	24958	54676	1.73%	0.182%
33	14.351	1105	1108	1112	rBV2	24256	55156	1.74%	0.184%
34	14.454	1115	1117	1120	rBV	22077	34972	1.10%	0.117%

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF120815\
Data File : BF083631.D
Acq On : 9 Dec 2015 9:39
Operator : UM/IZ
Sample : G4647-19
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
BRIDGE14-10(0-2)

Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF120815.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	14.557	1124	1126	1129	rBV	47572	73949	2.34%	0.247%
36	14.762	1140	1144	1147	rBV	39737	71439	2.26%	0.238%
37	14.865	1150	1153	1155	rBV	115756	155636	4.91%	0.519%
38	14.980	1161	1163	1167	rBV4	21605	55003	1.74%	0.183%
39	15.117	1173	1175	1178	rBV	116653	170437	5.38%	0.568%
40	15.414	1198	1201	1204	rBV	1255728	1757060	55.48%	5.858%
41	15.700	1223	1226	1230	rVB2	36524	64519	2.04%	0.215%
42	16.111	1260	1262	1266	rBV4	29774	54965	1.74%	0.183%
43	16.203	1266	1270	1271	rBV3	34760	91501	2.89%	0.305%
44	16.248	1272	1274	1276	rVB2	63746	88010	2.78%	0.293%
45	16.305	1276	1279	1284	rBV5	32830	72643	2.29%	0.242%
46	16.431	1288	1290	1292	rBV	24571	34650	1.09%	0.116%
47	16.774	1317	1320	1322	rBV	81098	109407	3.45%	0.365%
48	16.923	1331	1333	1337	rBV2	175836	358213	11.31%	1.194%
49	17.003	1337	1340	1343	rBV	516285	776667	24.52%	2.590%
50	17.357	1369	1371	1374	rBV3	45016	81858	2.58%	0.273%
51	17.746	1403	1405	1409	rBV	198520	263808	8.33%	0.880%
52	18.637	1482	1483	1489	rVB	286076	445557	14.07%	1.486%
53	21.872	1761	1766	1772	rVB	579128	1576616	49.79%	5.257%

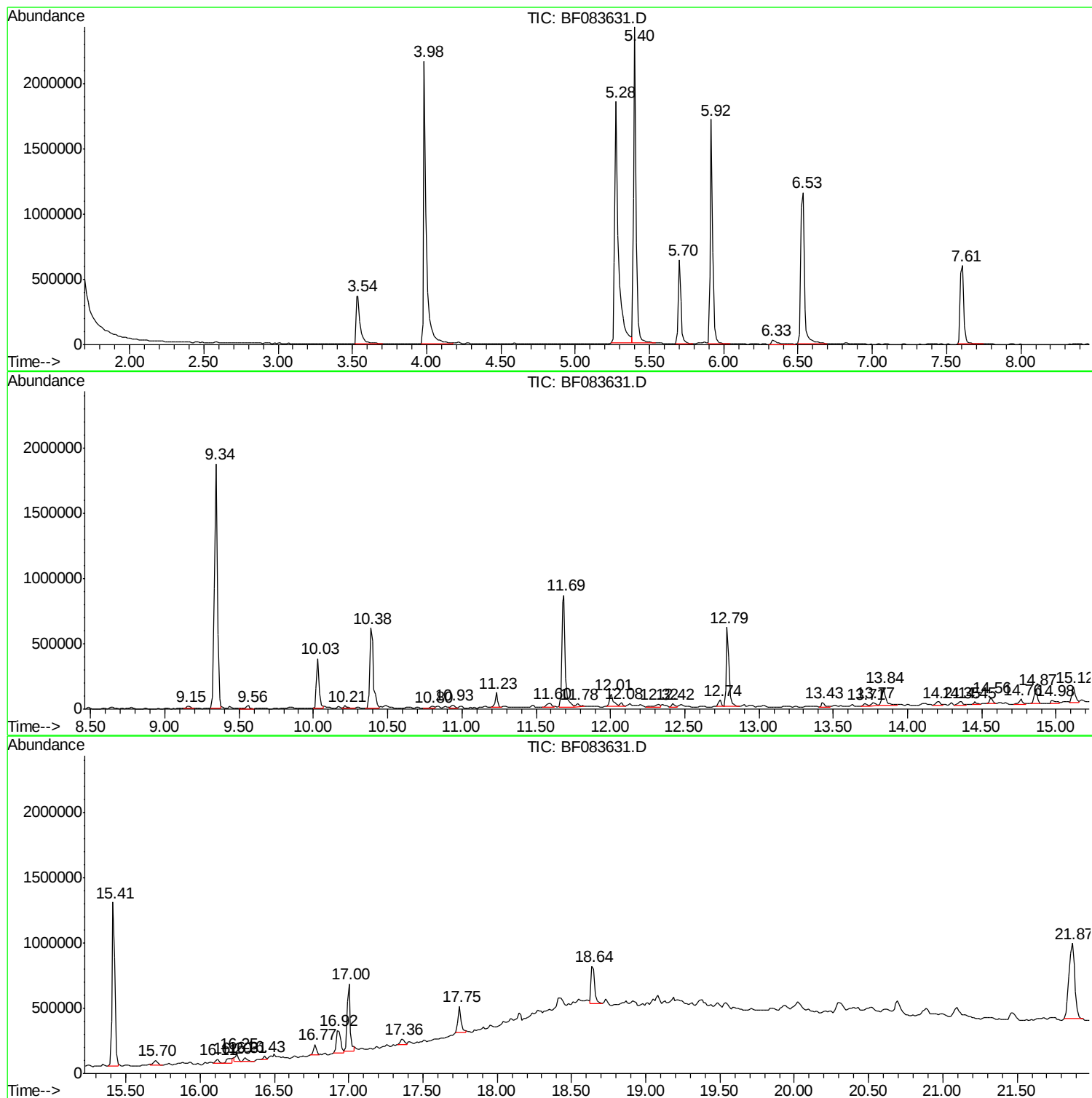
Sum of corrected areas: 29992707

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF120815\
Data File : BF083631.D
Acq On : 9 Dec 2015 9:39
Operator : UM/IZ
Sample : G4647-19
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
BRIDGE14-10(0-2)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF120815.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF120815\
Data File : BF083631.D
Acq On : 9 Dec 2015 9:39
Operator : UM/IZ
Sample : G4647-19
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
BRIDGE14-10(0-2)

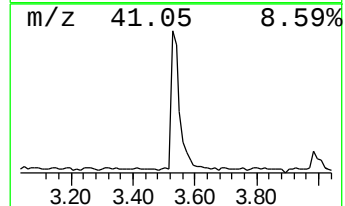
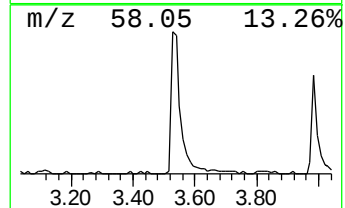
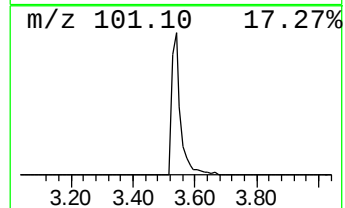
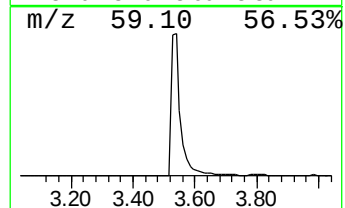
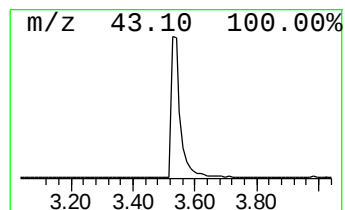
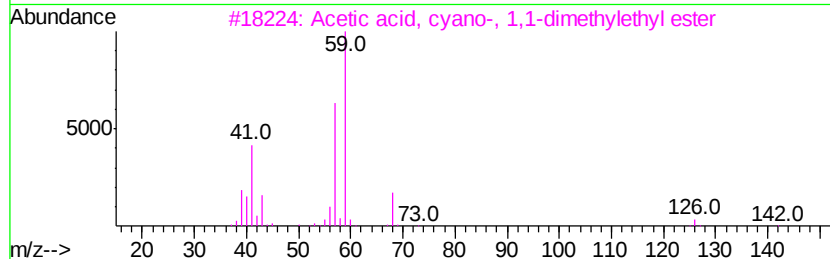
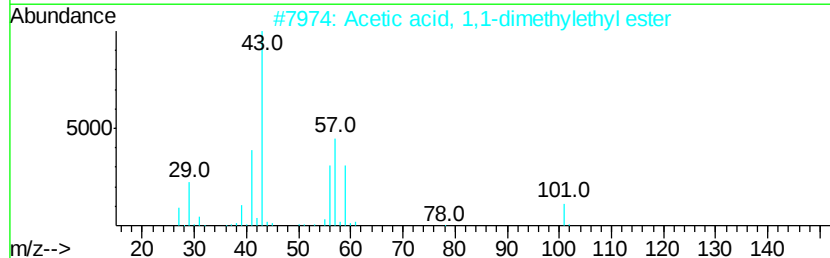
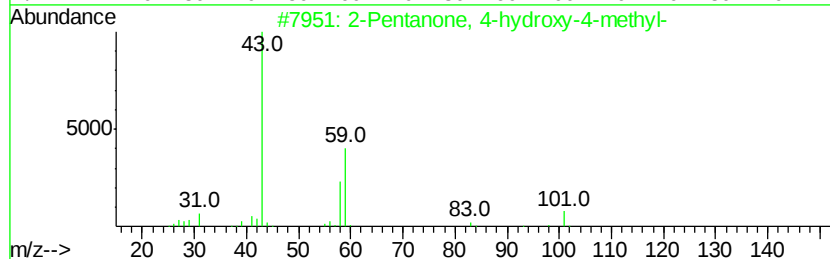
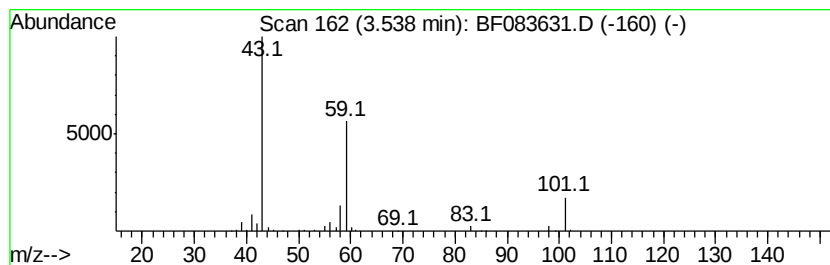
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF120815.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 unknown3.54 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.54	18.64 ng	753526	1,4-Dichlorobenzene-d4	5.70

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	45
2		Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	38
3		Acetic acid, cyano-, 1,1-dimethv...	141	C7H11NO2	001116-98-9	23
4		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
5		3-Hydroxy-3-methyl-2-butanone	102	C5H10O2	000115-22-0	9



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF120815\
Data File : BF083631.D
Acq On : 9 Dec 2015 9:39
Operator : UM/IZ
Sample : G4647-19
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
BRIDGE14-10(0-2)

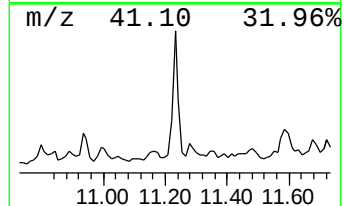
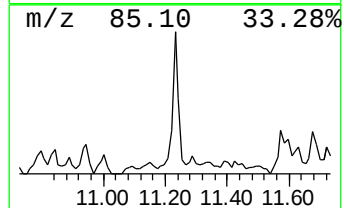
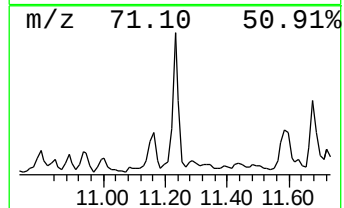
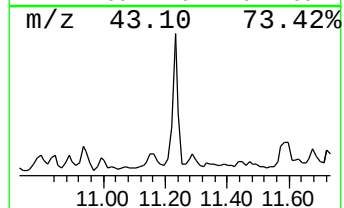
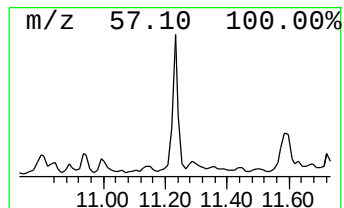
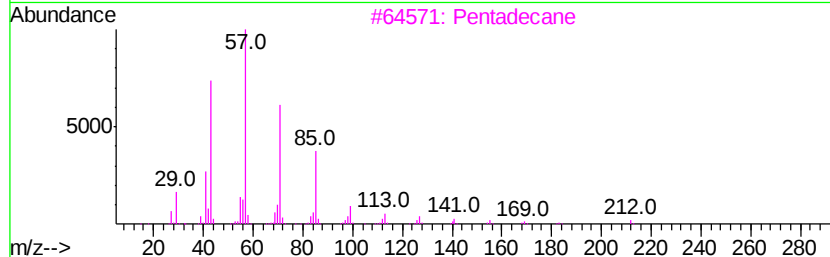
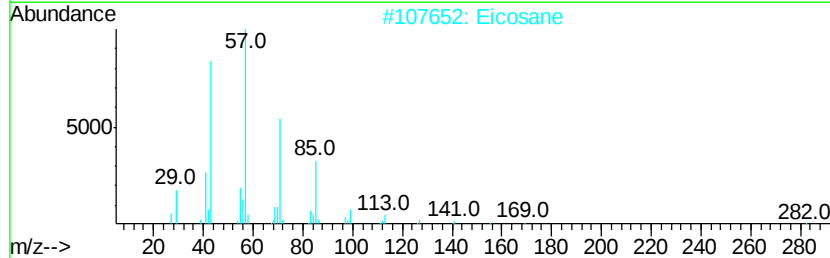
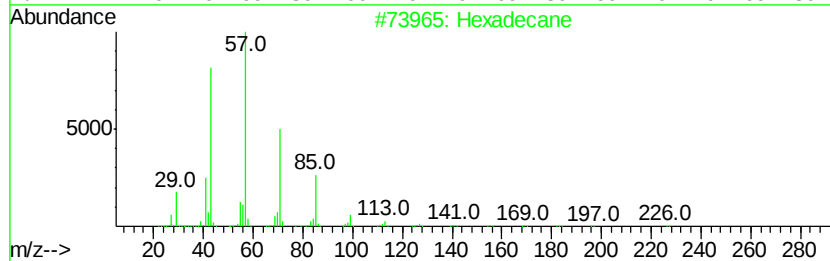
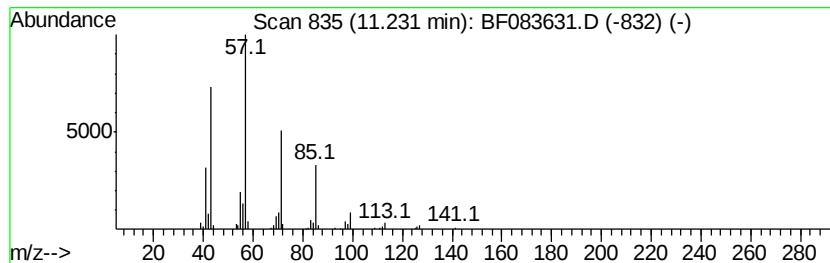
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF120815.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 4 Hexadecane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.23	2.79 ng	144458	Acenaphthene-d10	10.38

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecane	226	C16H34	000544-76-3	91
2		Eicosane	282	C20H42	000112-95-8	91
3		Pentadecane	212	C15H32	000629-62-9	90
4		Heptadecane	240	C17H36	000629-78-7	80
5		Tetradecane	198	C14H30	000629-59-4	72



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF120815\
Data File : BF083631.D
Acq On : 9 Dec 2015 9:39
Operator : UM/IZ
Sample : G4647-19
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
BRIDGE14-10(0-2)

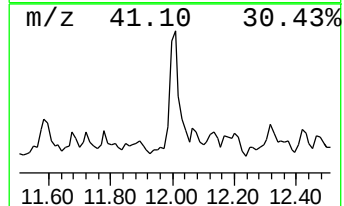
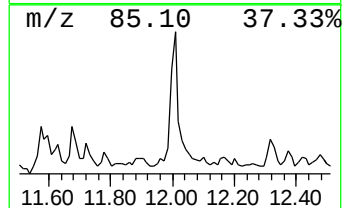
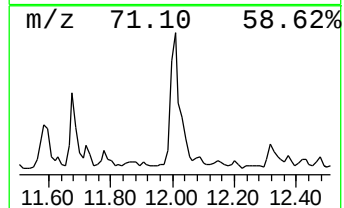
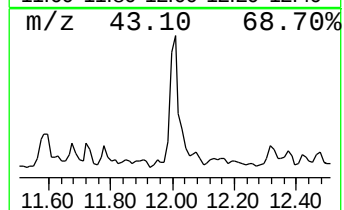
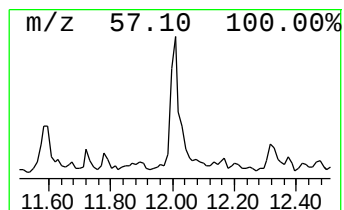
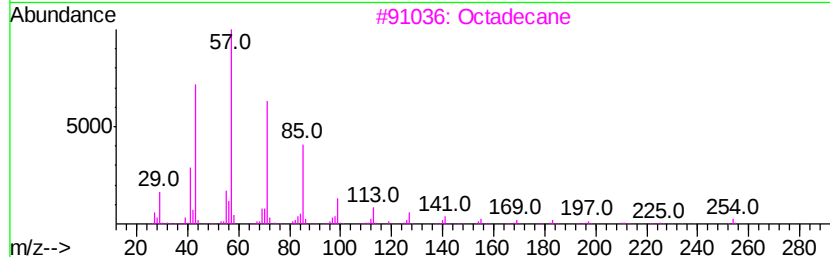
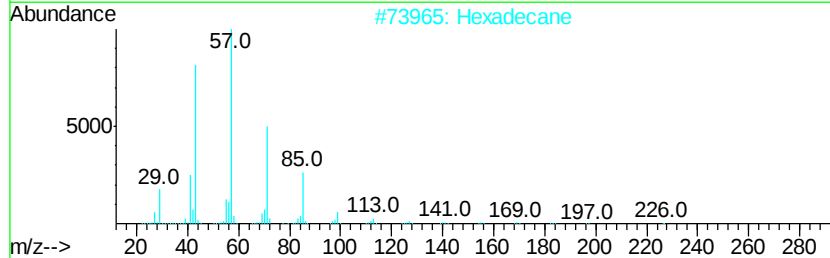
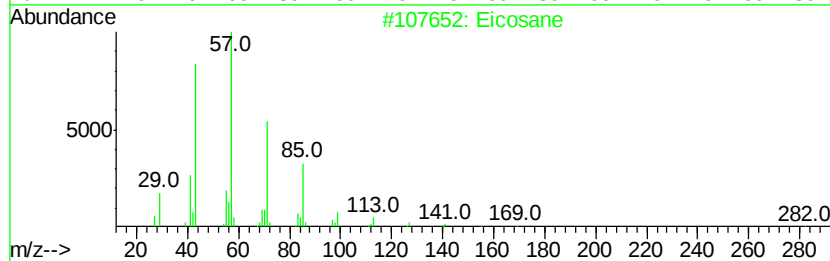
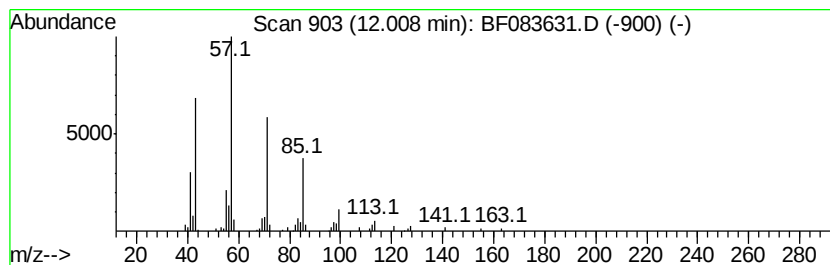
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF120815.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Eicosane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.01	4.66 ng	198004	Phenanthrene-d10	12.79

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	91
2		Hexadecane	226	C16H34	000544-76-3	90
3		Octadecane	254	C18H38	000593-45-3	78
4		Pentadecane	212	C15H32	000629-62-9	72
5		Heptacosane	380	C27H56	000593-49-7	64



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF120815\
Data File : BF083631.D
Acq On : 9 Dec 2015 9:39
Operator : UM/IZ
Sample : G4647-19
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
BRIDGE14-10(0-2)

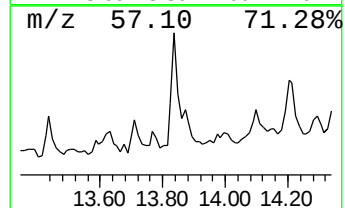
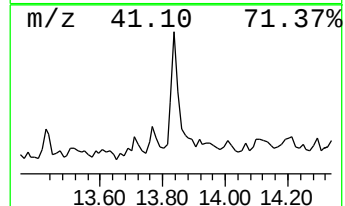
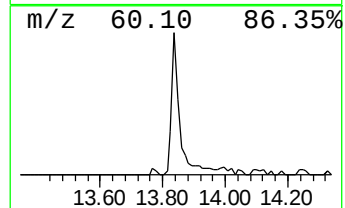
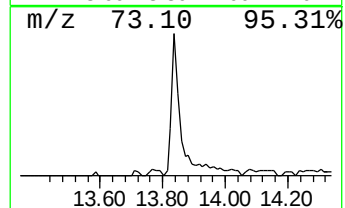
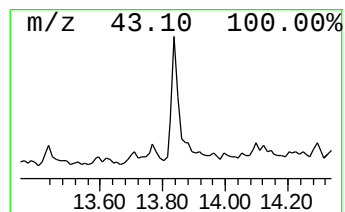
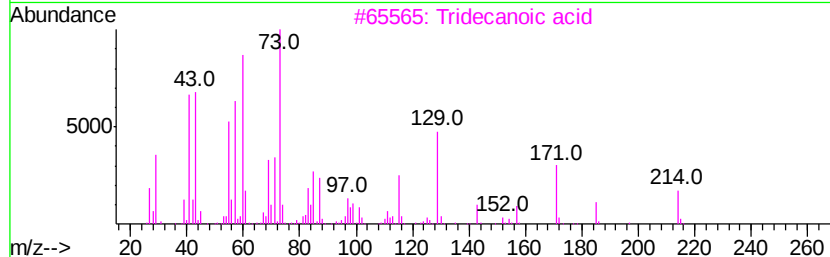
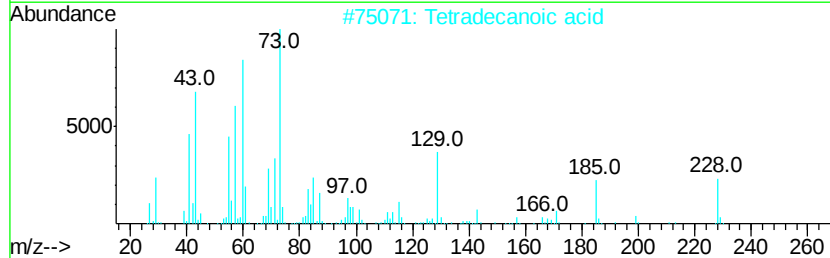
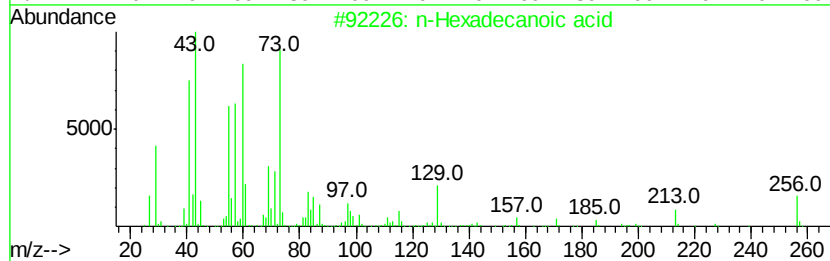
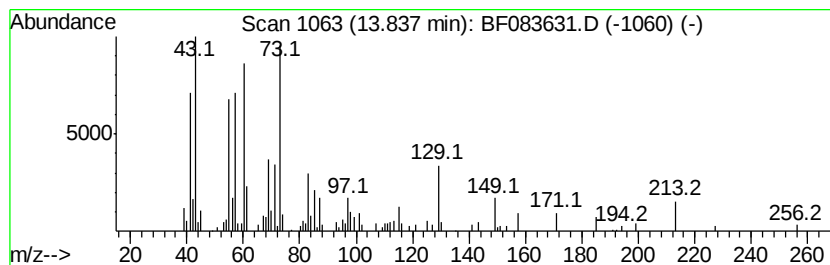
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF120815.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 n-Hexadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.84	5.37 ng	228160	Phenanthrene-d10	12.79

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	91
2		Tetradecanoic acid	228	C14H28O2	000544-63-8	72
3		Tridecanoic acid	214	C13H26O2	000638-53-9	64
4		Dodecanoic acid	200	C12H24O2	000143-07-7	64
5		Undecanoic acid	186	C11H22O2	000112-37-8	53



Data Path : Z:\HPCHEM1\BNA F\DATA\BF120815\
Data File : BF083631.D
Acq On : 9 Dec 2015 9:39
Operator : UM/IZ
Sample : G4647-19
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
BRIDGE14-10(0-2)

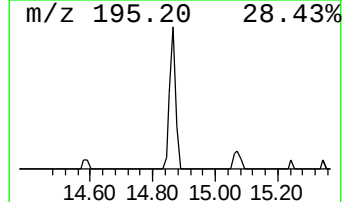
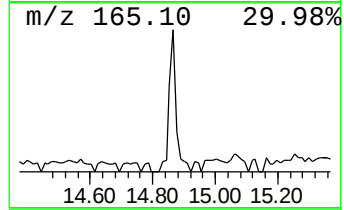
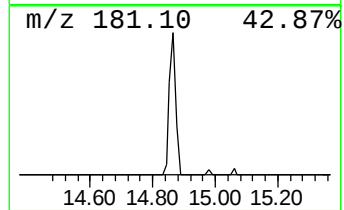
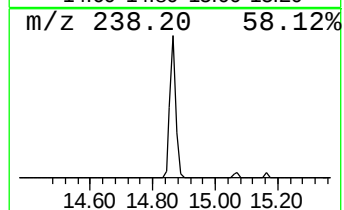
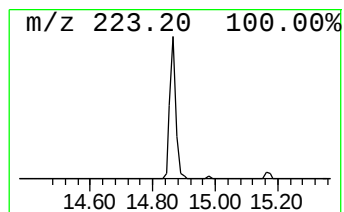
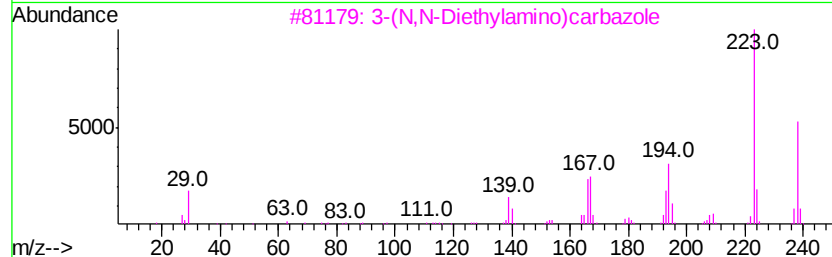
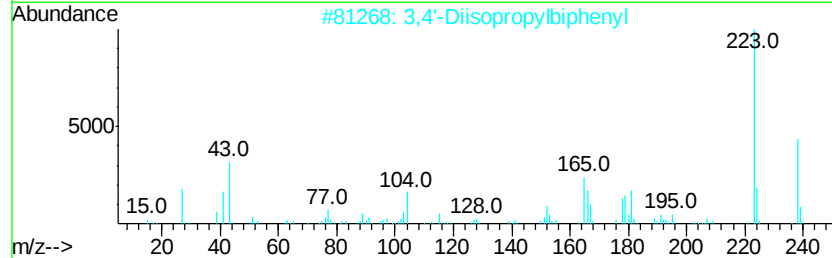
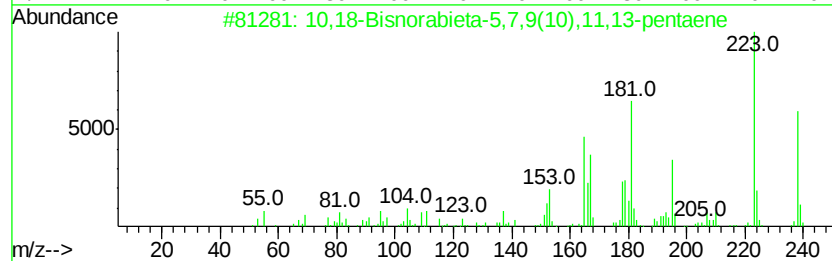
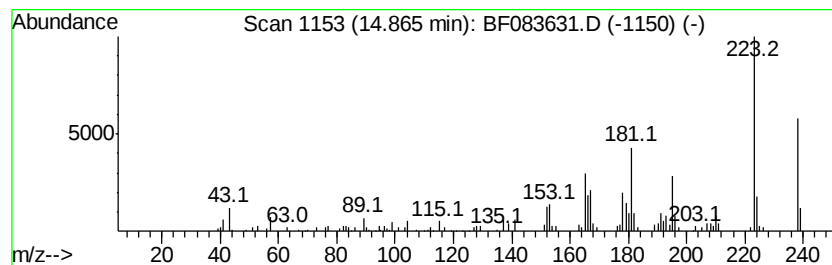
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF120815.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 10,18-Bisnorabieta-5,7,9(10... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.87	3.66 ng	155636	Phenanthrene-d10	12.79

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	10,18-Bisnorabieta-5,7,9(10),11,...	238	C18H22	006566-19-4	99
2		3,4'-Diisopropylbiphenyl	238	C18H22	061434-46-6	86
3		3-(N,N-Diethylamino)carbazole	238	C16H18N2	054994-31-9	53
4		4,4'-Diisopropylbiphenyl	238	C18H22	018970-30-4	53
5		Anthracene, 1,4-dimethoxy-	238	C16H14O2	013076-29-4	47



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF120815\
Data File : BF083631.D
Acq On : 9 Dec 2015 9:39
Operator : UM/IZ
Sample : G4647-19
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
BRIDGE14-10(0-2)

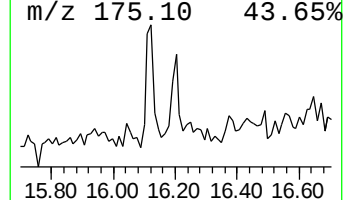
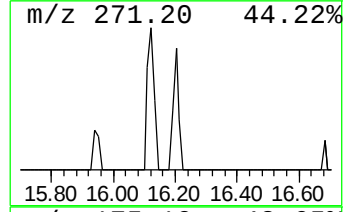
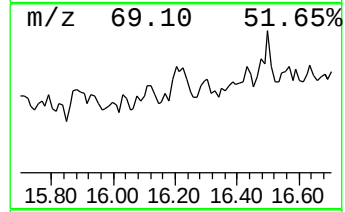
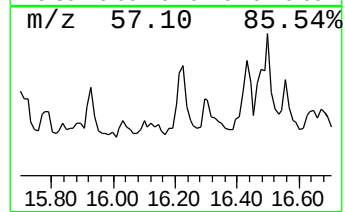
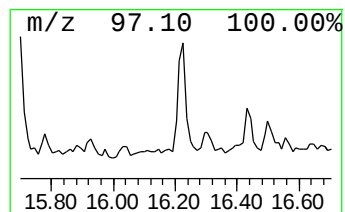
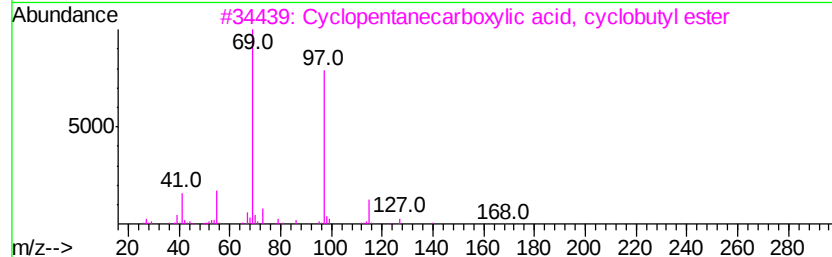
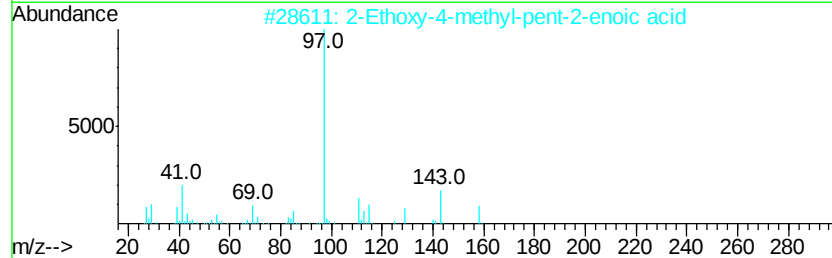
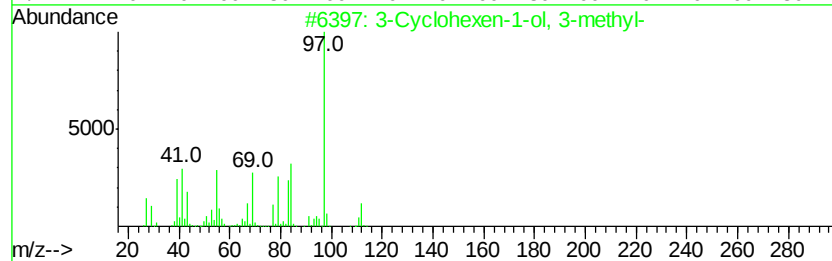
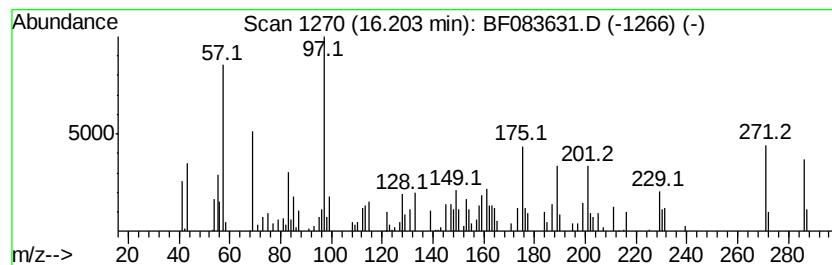
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF120815.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 unknown16.20 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.20	2.36 ng	91501	Chrysene-d12	17.00

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Cyclohexen-1-ol, 3-methyl-	112	C7H12O	053783-91-8	14
2		2-Ethoxy-4-methyl-pent-2-enoic acid	158	C8H14O3	1000190-08-8	14
3		Cyclopentanecarboxylic acid, cyclobutyl ester	168	C10H16O2	1000282-76-8	14
4		4,5-Dihydro-N-hydroxy-N-phenyl-3-pyridinecarboxamide	205	C11H11N03	1000243-32-3	14
5		Cyclohexane, 1,1-dimethyl-	112	C8H16	000590-66-9	14



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF120815\
Data File : BF083631.D
Acq On : 9 Dec 2015 9:39
Operator : UM/IZ
Sample : G4647-19
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
BRIDGE14-10(0-2)

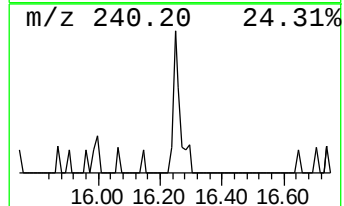
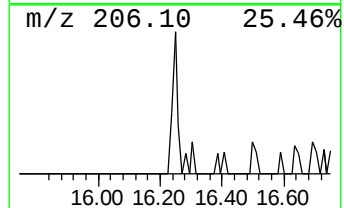
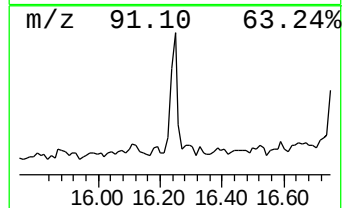
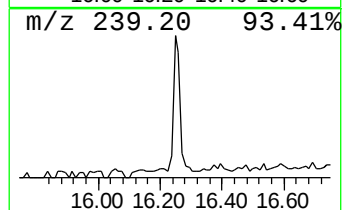
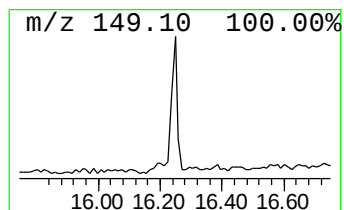
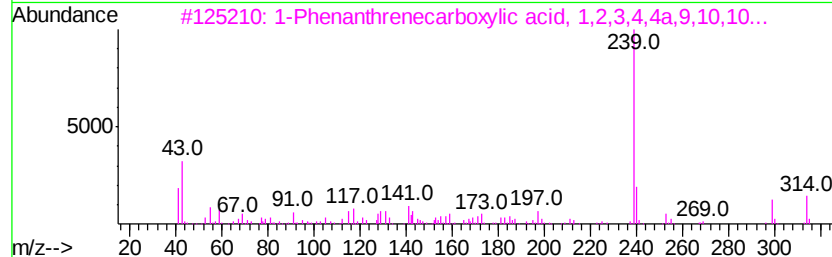
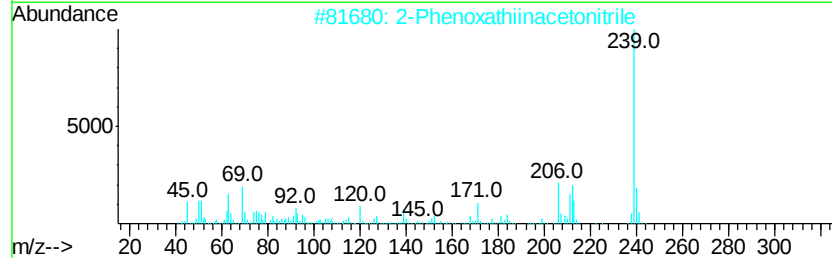
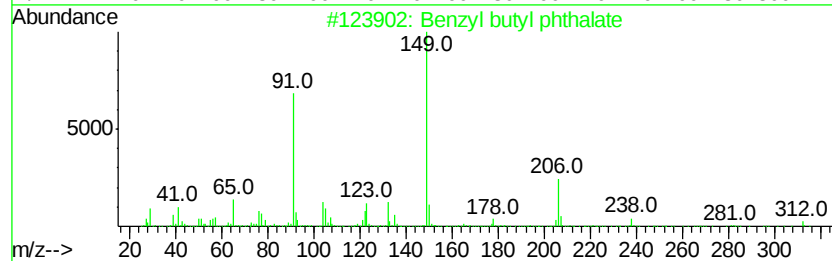
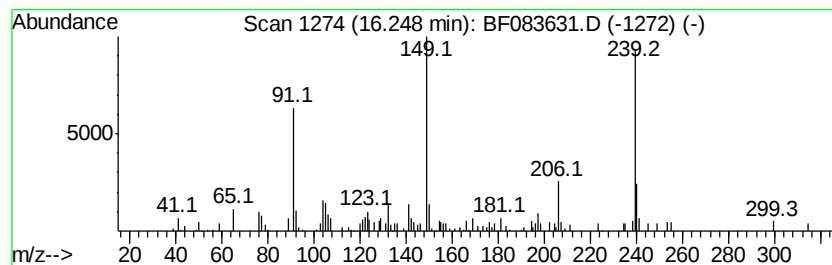
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF120815.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 unknown16.25 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.25	2.27 ng	88010	Chrysene-d12	17.00

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzyl butyl phthalate	312	C19H20O4	000085-68-7	41
2		2-Phenoxathiinacetonitrile	239	C14H9NOS	1000255-56-5	38
3		1-Phenanthrenecarboxylic acid, 1...	314	C21H30O2	001235-74-1	30
4		4-Amino-3-(N,N-dimethylamino)-9-...	239	C15H17N3	094127-27-2	27
5		9,10-Anthracenedione, 1-amino-4-...	239	C14H9NO3	000116-85-8	27



Data Path : Z:\HPCHEM1\BNA F\DATA\BF120815\
Data File : BF083631.D
Acq On : 9 Dec 2015 9:39
Operator : UM/IZ
Sample : G4647-19
Misc :
ALS Vial : 39 Sample Multiplier: 1

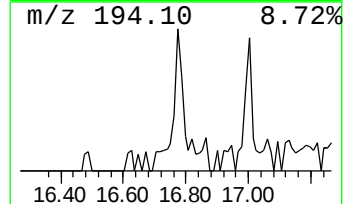
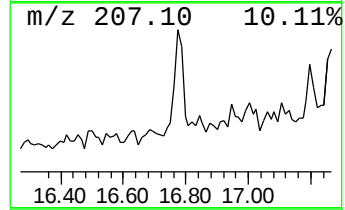
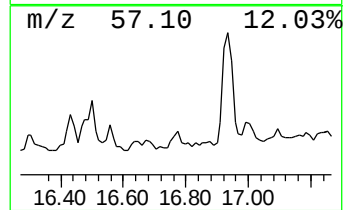
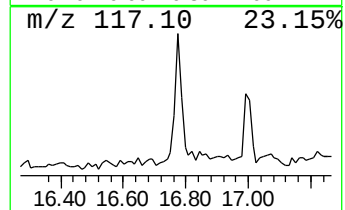
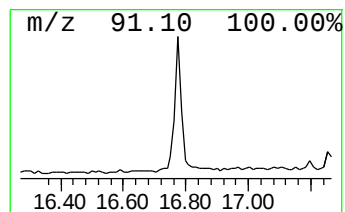
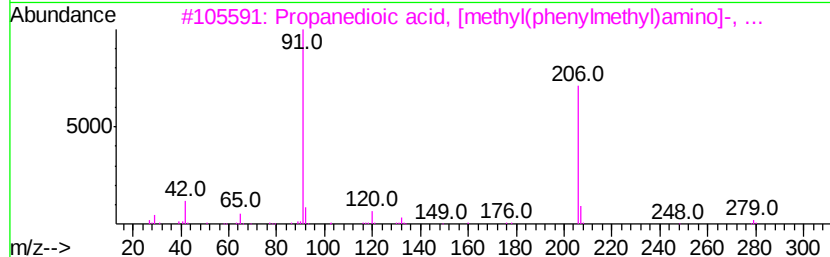
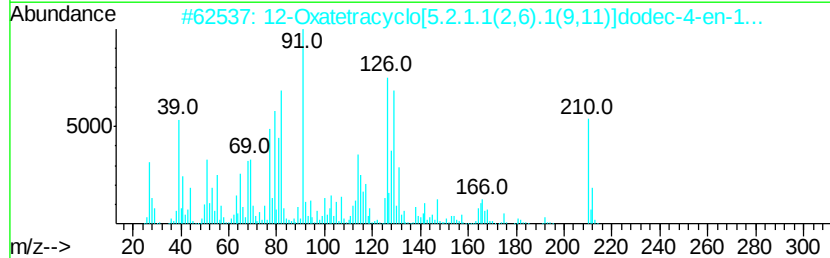
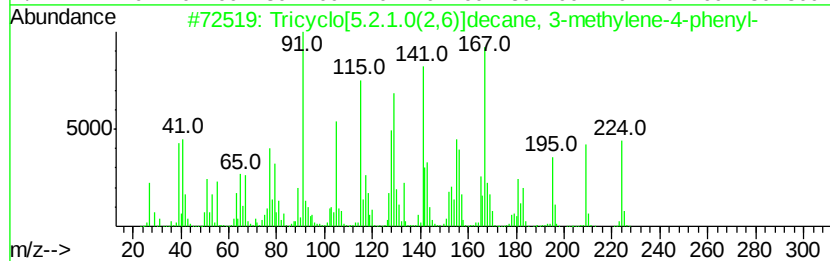
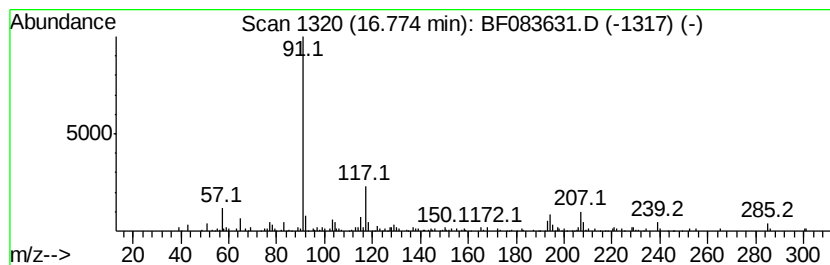
Instrument :
BNA_F
ClientSampleId :
BRIDGE14-10(0-2)

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF120815.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Tricyclo[5.2.1.0(2,6)]decan... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.		
16.77	2.82 ng	109407	Chrysene-d12	17.00		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tricyclo[5.2.1.0(2,6)]decane, 3-...	224	C17H20	1000150-37-1	56
2		12-Oxatetracyclo[5.2.1.1(2,6).1(...	210	C11H11ClO2	078630-89-4	38
3		Propanedioic acid, [methvl(phenyl...	279	C15H21N04	001032-02-6	27
4		Cyclohexanone, 2-[bis(phenylmet...	383	C27H29NO	094318-25-9	16
5		Benzonitrile, m-phenethyl-	207	C15H13N	034176-91-5	16



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF120815\
Data File : BF083631.D
Acq On : 9 Dec 2015 9:39
Operator : UM/IZ
Sample : G4647-19
Misc :
ALS Vial : 39 Sample Multiplier: 1

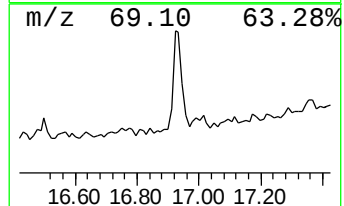
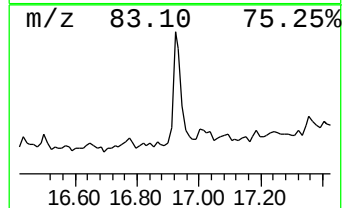
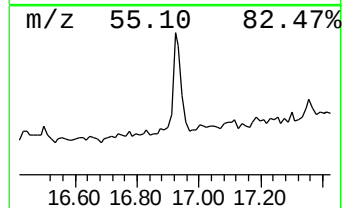
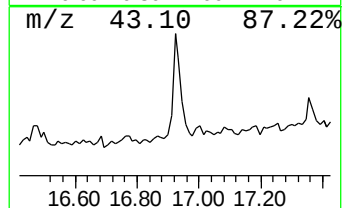
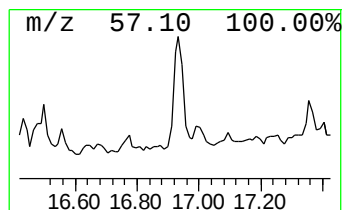
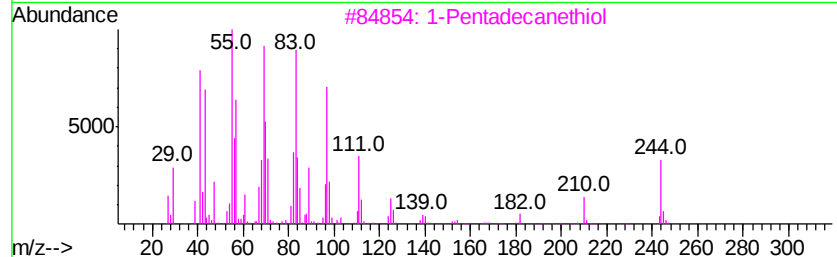
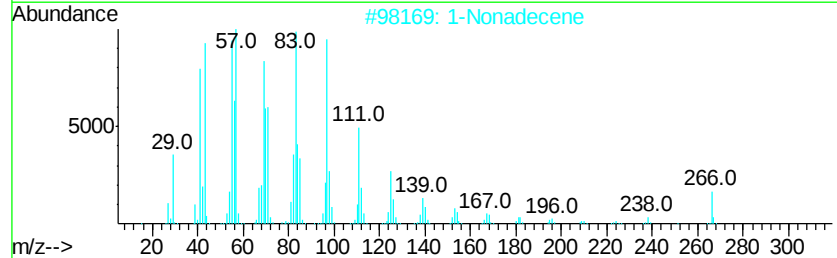
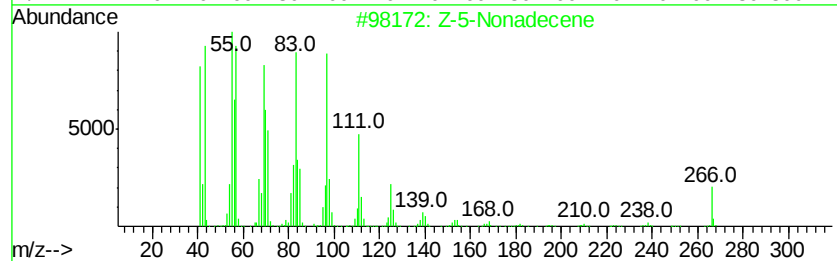
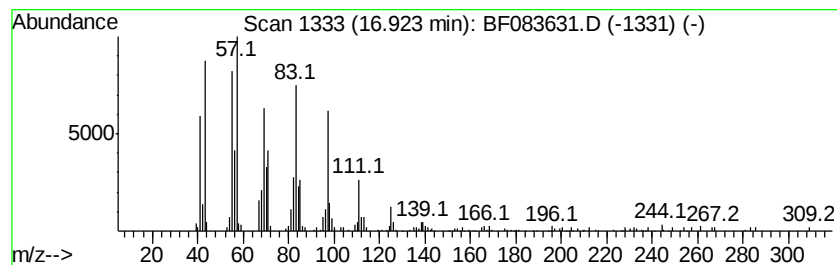
Instrument :
BNA_F
ClientSampleId :
BRIDGE14-10(0-2)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF120815.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Z-5-Nonadecene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD		R.T.
16.92	9.22 ng	358213	Chrysene-d12		17.00
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Z-5-Nonadecene	266	C19H38	1000131-11-8	98
2	1-Nonadecene	266	C19H38	018435-45-5	97
3	1-Pentadecanethiol	244	C15H32S	025276-70-4	95
4	5-Eicosene, (E)-	280	C20H40	074685-30-6	91
5	1-Docosene	308	C22H44	001599-67-3	91



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF120815\
Data File : BF083631.D
Acq On : 9 Dec 2015 9:39
Operator : UM/IZ
Sample : G4647-19
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
BRIDGE14-10(0-2)

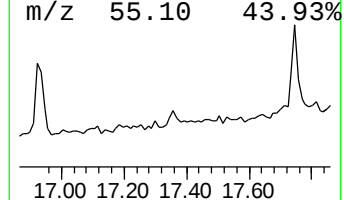
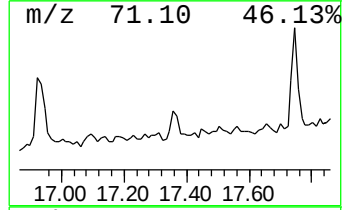
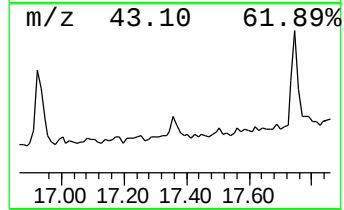
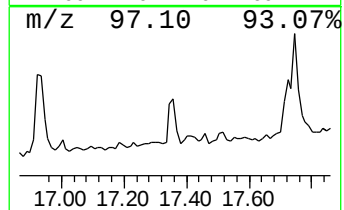
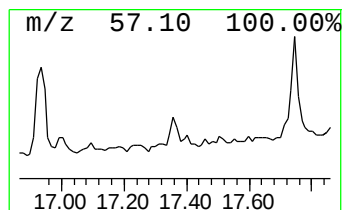
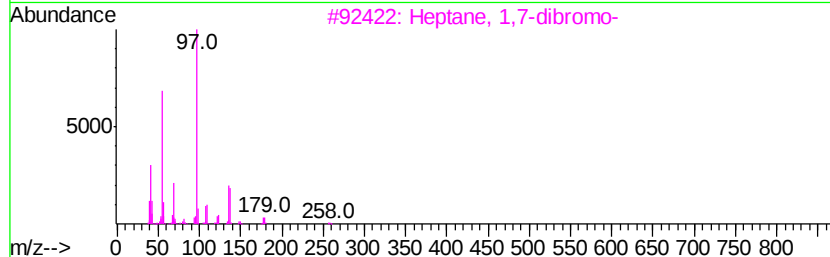
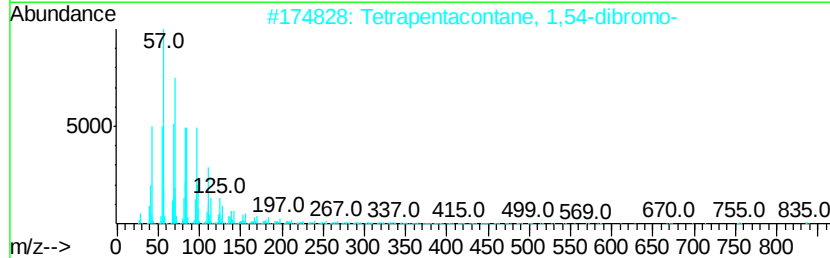
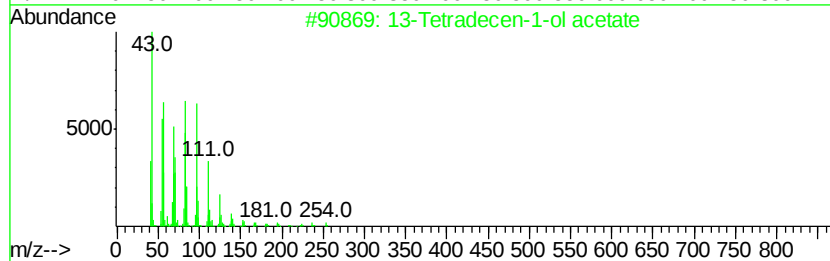
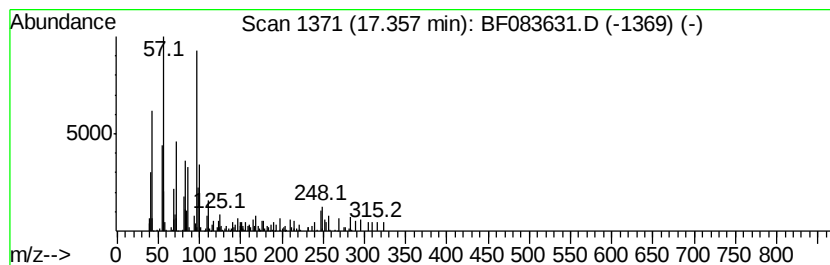
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF120815.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 13-Tetradecen-1-ol acetate Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.36	2.11 ng	81858	Chrysene-d12	17.00

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			13-Tetradecen-1-ol acetate	254	C16H30O2	056221-91-1	59
2			Tetrapentacontane, 1,54-dibromo-	915	C54H108Br2	1000156-09-4	27
3			Heptane, 1,7-dibromo-	256	C7H14Br2	004549-31-9	27
4			2-Propionyl-6-methyl-3,4-dihydro...	154	C9H14O2	1000145-07-7	27
5			2,4:3,5-Diethylidene-1-xylose	202	C9H14O5	1000126-95-9	25



Data Path : Z:\HPCHEM1\BNA F\DATA\BF120815\
Data File : BF083631.D
Acq On : 9 Dec 2015 9:39
Operator : UM/IZ
Sample : G4647-19
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
BRIDGE14-10(0-2)

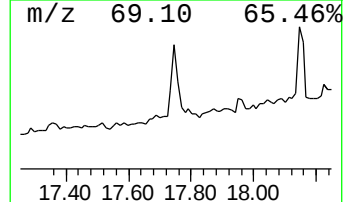
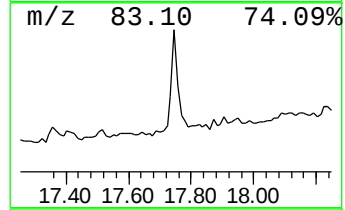
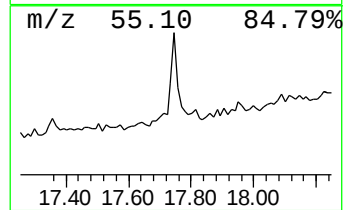
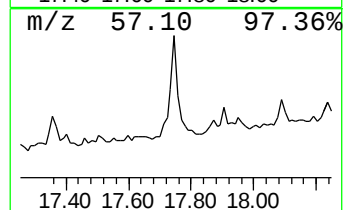
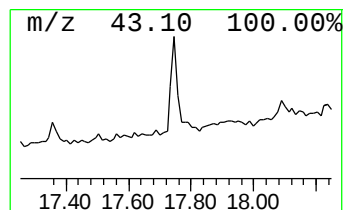
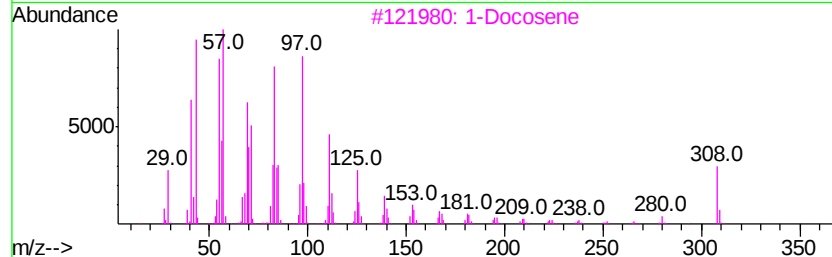
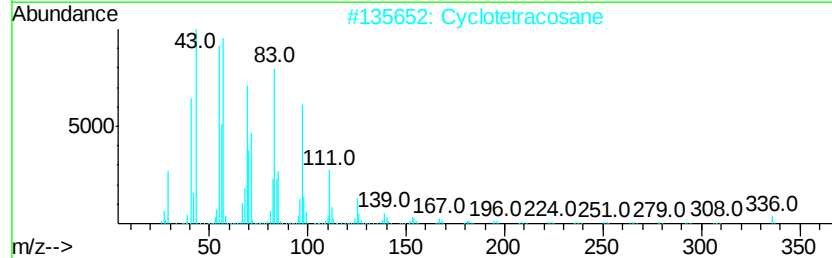
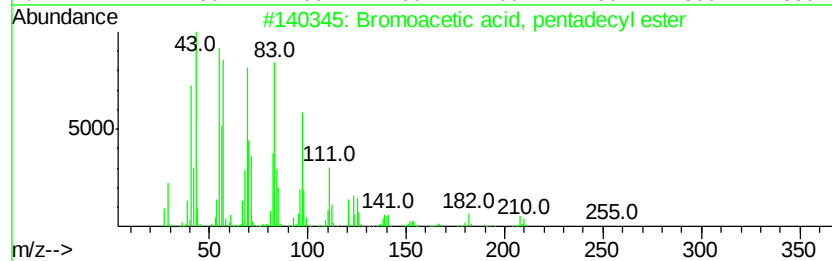
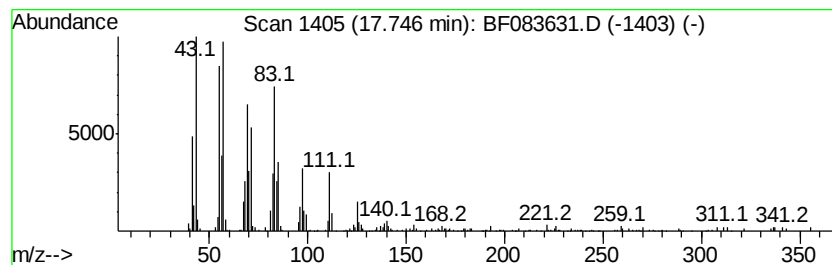
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF120815.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Bromoacetic acid, pentadecyl... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.75	6.79 ng	263808	Chrysene-d12	17.00

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Bromoacetic acid, pentadecyl ester	348	C17H33BrO2	131143-01-6	74
2			Cyclotetracosane	336	C24H48	000297-03-0	68
3			1-Docosene	308	C22H44	001599-67-3	68
4			Trifluoroacetic acid, n-octadecyl...	366	C20H37F3O2	079392-43-1	58
5			Cyclopentane, 1-ethyl-3-methyl-,...	112	C8H16	002613-66-3	53



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF120815\
Data File : BF083631.D
Acq On : 9 Dec 2015 9:39
Operator : UM/IZ
Sample : G4647-19
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
BRIDGE14-10(0-2)

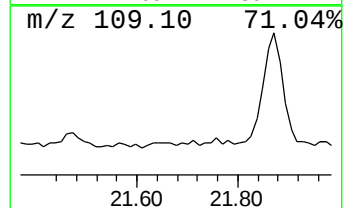
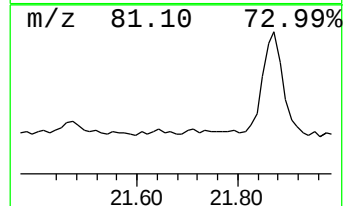
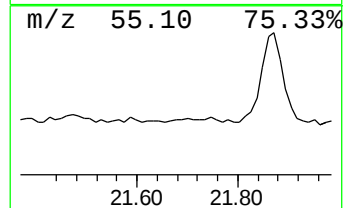
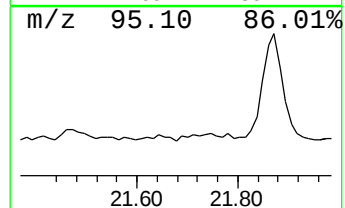
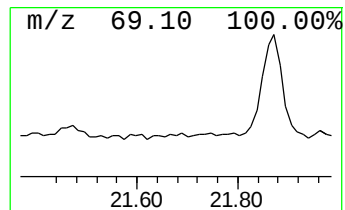
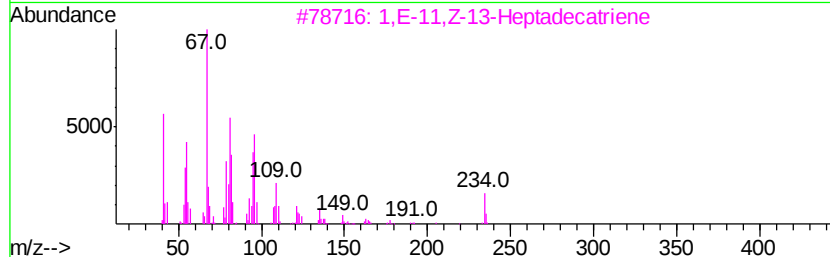
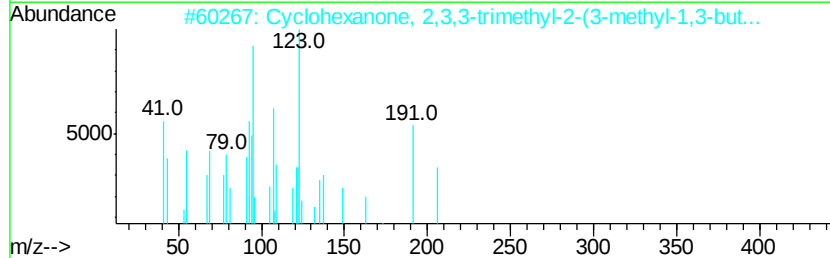
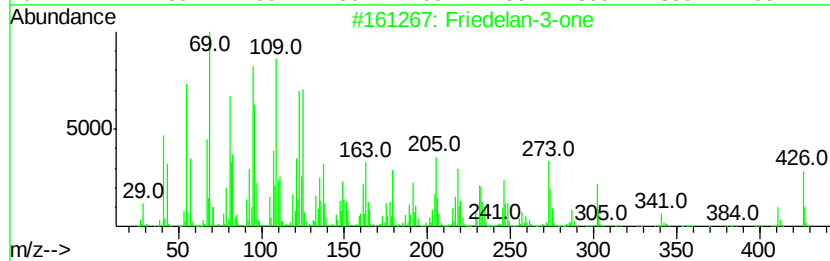
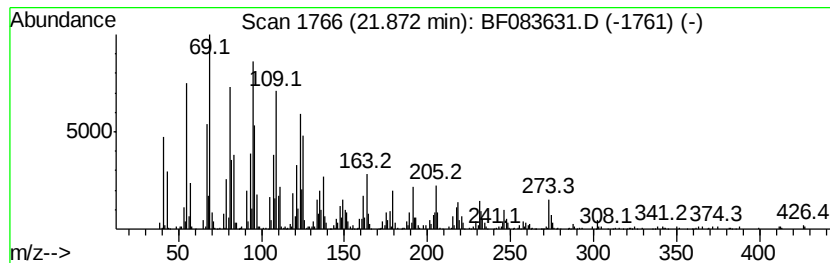
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF120815.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Friedelan-3-one Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.87	70.77 ng	1576620	Perylene-d12	18.64

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Friedelan-3-one	426	C30H50O	000559-74-0	95
2		Cyclohexanone, 2,3,3-trimethyl-2...	206	C14H22O	069296-90-8	52
3		1,E-11,Z-13-Heptadecatriene	234	C17H30	080625-35-0	44
4		2,3,3-Trimethyl-2-(3-methyl-buta...	206	C14H22O	1000193-60-6	38
5		3-Cyclopentylpropionic acid, but...	194	C12H18O2	1000292-46-4	30



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF120815\
 Data File : BF083631.D
 Acq On : 9 Dec 2015 9:39
 Operator : UM/IZ
 Sample : G4647-19
 Misc :
 ALS Vial : 39 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 BRIDGE14-10(0-2)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF120815.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
unknown3.54	3.54	18.6	ng	753526	1	5.70	808713	20.0
Hexadecane	11.23	2.8	ng	144458	3	10.38	1034300	20.0
Eicosane	12.01	4.7	ng	198004	4	12.79	849923	20.0
n-Hexadecanoic acid	13.84	5.4	ng	228160	4	12.79	849923	20.0
10,18-Bisnorabiet...	14.87	3.7	ng	155636	4	12.79	849923	20.0
unknown16.20	16.20	2.4	ng	91501	5	17.00	776667	20.0
unknown16.25	16.25	2.3	ng	88010	5	17.00	776667	20.0
Tricyclo[5.2.1.0(...	16.77	2.8	ng	109407	5	17.00	776667	20.0
Z-5-Nonadecene	16.92	9.2	ng	358213	5	17.00	776667	20.0
13-Tetradecen-1-o...	17.36	2.1	ng	81858	5	17.00	776667	20.0
Bromoacetic acid,...	17.75	6.8	ng	263808	5	17.00	776667	20.0
Friedelan-3-one	21.87	70.8	ng	1576620	6	18.64	445557	20.0