

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF120815\
 Data File : BF083624.D
 Acq On : 9 Dec 2015 6:08
 Operator : UM/IZ
 Sample : G4647-05
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 BRIDGE14-3(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	1.847	13	14	34	rVB3	114545	478613	16.18%	1.672%
2	3.527	159	161	173	rBV	373951	681693	23.05%	2.382%
3	3.984	199	201	219	rBV	1726537	2743177	92.76%	9.586%
4	5.264	311	313	322	rBV	1462852	2885958	97.59%	10.085%
5	5.402	322	325	344	rVB	1762665	2957175	100.00%	10.333%
6	5.699	349	351	359	rBV	738189	933050	31.55%	3.260%
7	5.916	367	370	382	rVB	1452061	1939344	65.58%	6.777%
8	6.339	405	407	415	rBV	18887	45504	1.54%	0.159%
9	6.522	421	423	438	rBV	1183427	1803552	60.99%	6.302%
10	7.596	514	517	529	rBV	853881	1153385	39.00%	4.030%
11	9.345	667	670	677	rBV	1616120	2431522	82.22%	8.497%
12	10.031	727	730	735	rVB	328788	472347	15.97%	1.651%
13	10.385	758	761	768	rBV	848757	1175515	39.75%	4.108%
14	10.796	790	797	799	rBV3	13077	36098	1.22%	0.126%
15	10.934	806	809	812	rVB	22960	33965	1.15%	0.119%
16	11.151	821	828	832	rBV2	10196	31338	1.06%	0.110%
17	11.231	832	835	838	rVV	94643	124190	4.20%	0.434%
18	11.585	862	866	868	rBV	30211	64121	2.17%	0.224%
19	11.676	871	874	881	rVV	872834	1244580	42.09%	4.349%
20	11.997	899	902	906	rBV	95896	167112	5.65%	0.584%
21	12.065	906	908	911	rVV2	22502	38885	1.31%	0.136%
22	12.339	931	932	937	rVB4	15557	37352	1.26%	0.131%
23	12.419	937	939	941	rBV	20517	31136	1.05%	0.109%
24	12.728	964	966	968	rBV	39874	51731	1.75%	0.181%
25	12.785	968	971	974	rVV	625996	896461	30.31%	3.133%
26	13.425	1024	1027	1030	rBV4	22457	30402	1.03%	0.106%
27	13.837	1060	1063	1069	rBV2	95888	183410	6.20%	0.641%
28	14.202	1092	1095	1097	rVV	20269	34676	1.17%	0.121%
29	14.351	1105	1108	1112	rBV2	32215	57905	1.96%	0.202%
30	14.557	1124	1126	1129	rBV	42511	65787	2.22%	0.230%
31	14.762	1141	1144	1148	rVB	28144	44121	1.49%	0.154%
32	14.854	1149	1152	1155	rBV	86594	113329	3.83%	0.396%
33	14.968	1159	1162	1167	rBV4	42258	107924	3.65%	0.377%
34	15.105	1172	1174	1178	rBV	90312	165423	5.59%	0.578%

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF120815\
Data File : BF083624.D
Acq On : 9 Dec 2015 6:08
Operator : UM/IZ
Sample : G4647-05
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
BRIDGE14-3(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	15.414	1197	1201	1206	rVB	912166	1450032	49.03%	5.067%
36	15.654	1219	1222	1223	rBV3	20308	36204	1.22%	0.127%
37	15.688	1223	1225	1229	rVB3	28626	56913	1.92%	0.199%
38	16.100	1259	1261	1265	rVB5	28928	63765	2.16%	0.223%
39	16.180	1265	1268	1269	rBV3	35002	65754	2.22%	0.230%
40	16.237	1271	1273	1276	rVV3	41607	84573	2.86%	0.296%
41	16.294	1276	1278	1281	rVV4	24772	46482	1.57%	0.162%
42	16.488	1293	1295	1299	rBV4	24757	62475	2.11%	0.218%
43	16.626	1305	1307	1309	rBV3	29239	47622	1.61%	0.166%
44	16.763	1317	1319	1322	rVB	45580	78963	2.67%	0.276%
45	16.923	1330	1333	1336	rBV2	161491	319369	10.80%	1.116%
46	16.991	1336	1339	1346	rVV	660604	971545	32.85%	3.395%
47	17.734	1402	1404	1412	rBV	182292	296565	10.03%	1.036%
48	18.409	1461	1463	1466	rBV2	141651	207641	7.02%	0.726%
49	18.637	1481	1483	1489	rVB	437776	633247	21.41%	2.213%
50	20.866	1675	1678	1682	rVB2	172164	371943	12.58%	1.300%
51	21.837	1760	1763	1769	rVB2	204464	563539	19.06%	1.969%

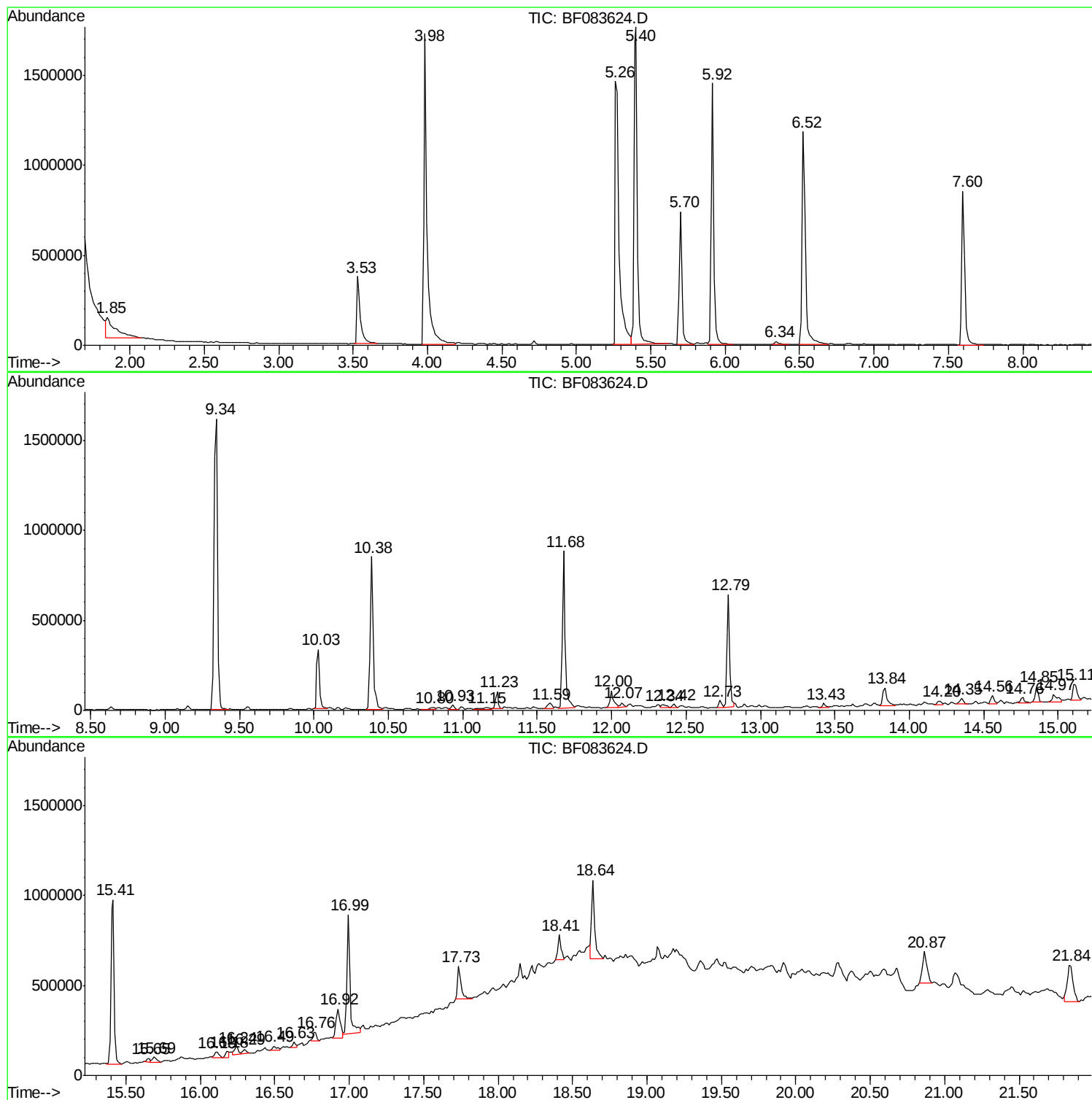
Sum of corrected areas: 28617413

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF120815\
Data File : BF083624.D
Acq On : 9 Dec 2015 6:08
Operator : UM/IZ
Sample : G4647-05
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
BRIDGE14-3(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 : BF083624.D
Acq On : 9 Dec 2015 6:08
Operator : UM/IZ
Sample : G4647-05
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
BRIDGE14-3(0-2)

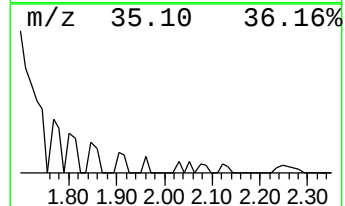
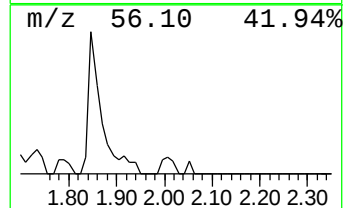
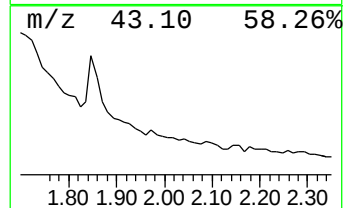
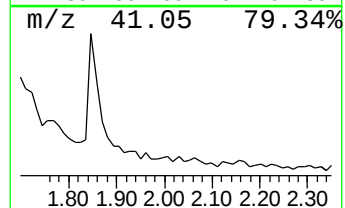
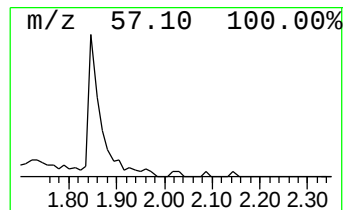
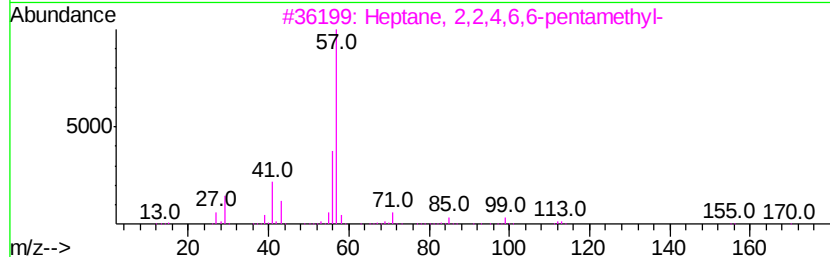
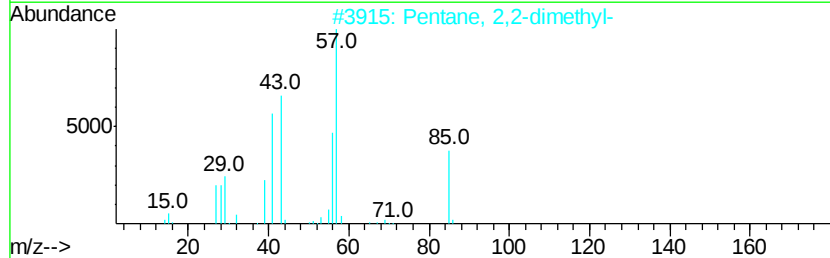
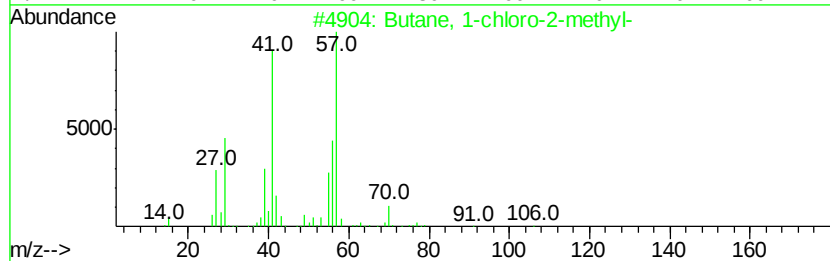
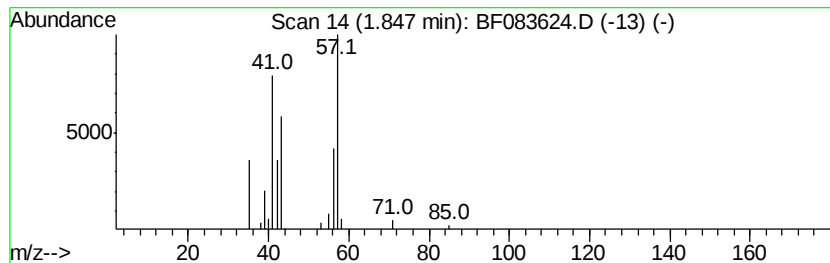
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF120815.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 unknown1.85 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.85	10.26 ng	478613	1,4-Dichlorobenzene-d4	5.70

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 1-chloro-2-methyl-	106	C5H11Cl	000616-13-7	32
2		Pentane, 2,2-dimethyl-	100	C7H16	000590-35-2	16
3		Heptane, 2,2,4,6,6-pentamethyl-	170	C12H26	013475-82-6	12
4		Pentane, 2,4-dimethyl-	100	C7H16	000108-08-7	10
5		2-Butene, (E)-	56	C4H8	000624-64-6	9



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF120815\
Data File : BF083624.D
Acq On : 9 Dec 2015 6:08
Operator : UM/IZ
Sample : G4647-05
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
BRIDGE14-3(0-2)

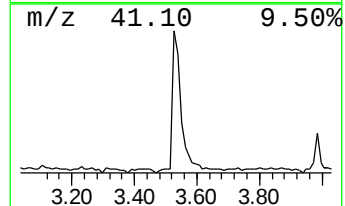
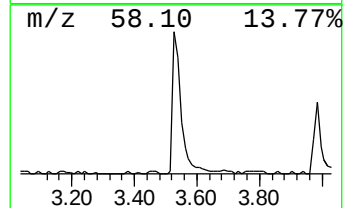
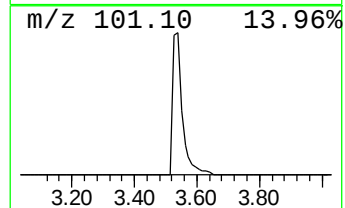
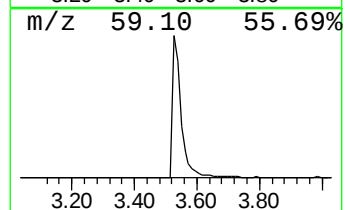
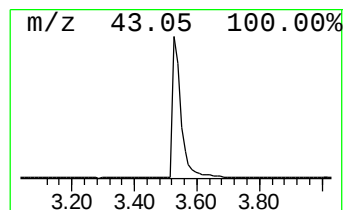
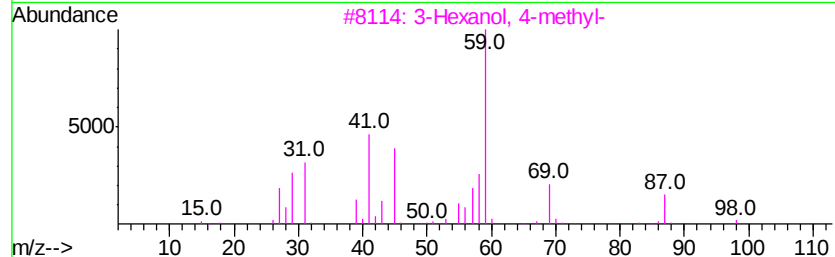
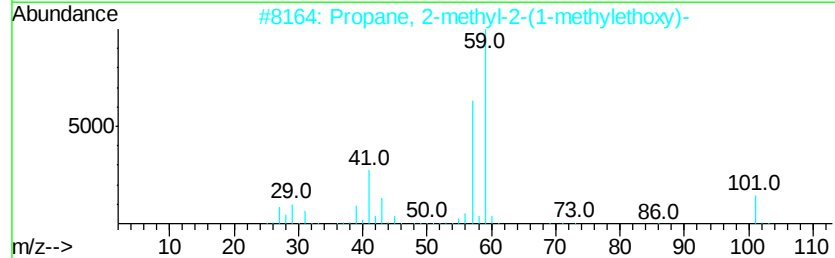
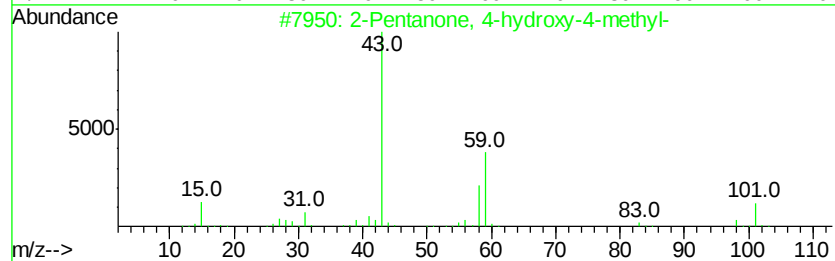
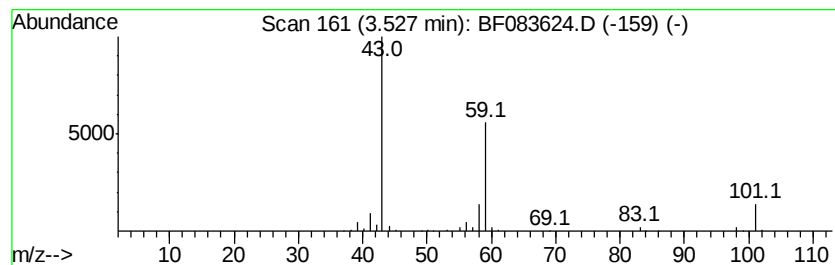
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF120815.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.53	14.61 ng	681693	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	50
2		Propane, 2-methyl-2-(1-methylethoxy)-	116	C7H16O	017348-59-3	42
3		3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
4		2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	33
5		Acetic acid, cyano-, 1,1-dimethyl-	141	C7H11NO2	001116-98-9	23



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF120815\
Data File : BF083624.D
Acq On : 9 Dec 2015 6:08
Operator : UM/IZ
Sample : G4647-05
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
BRIDGE14-3(0-2)

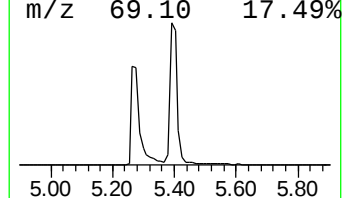
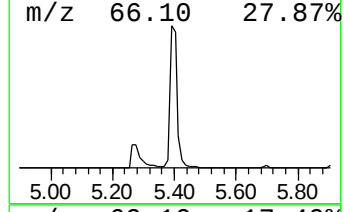
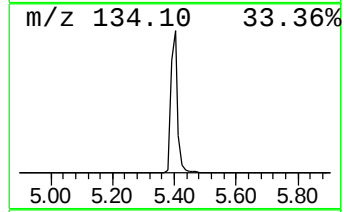
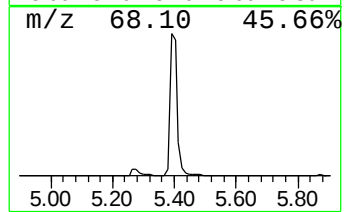
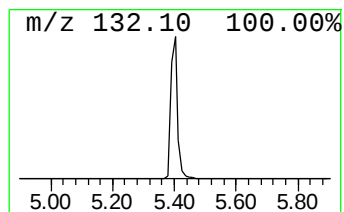
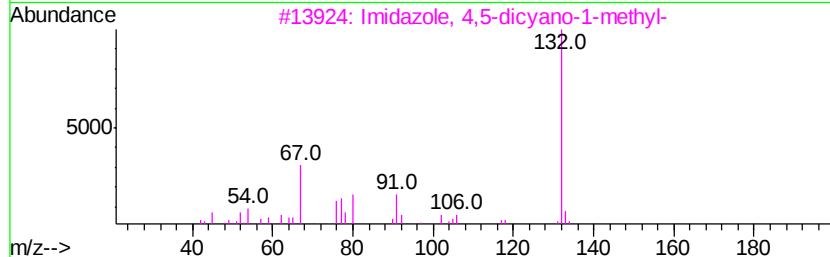
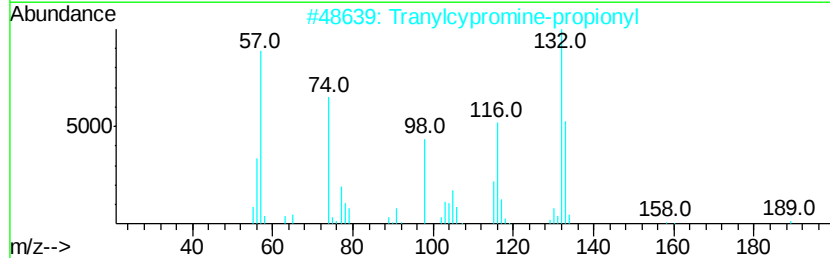
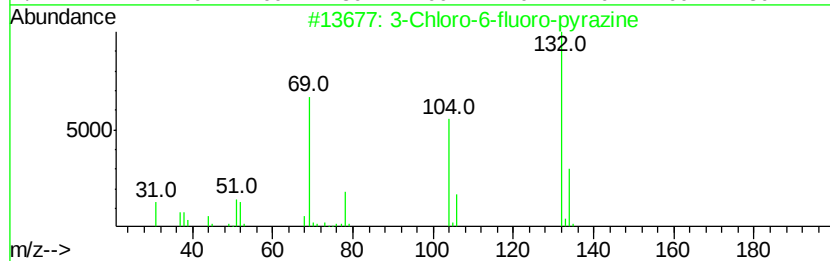
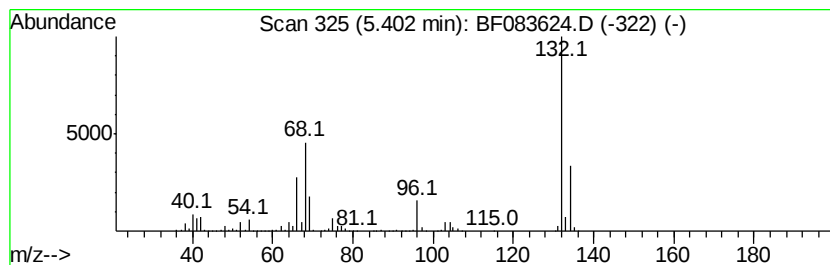
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF120815.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 unknown5.40 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.40	63.39 ng	2957180	1,4-Dichlorobenzene-d4	5.70

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
3		Imidazole, 4,5-dicyano-1-methyl-	132	C6H4N4	019485-35-9	9
4		Benzene, 1,3,5-trifluoro-	132	C6H3F3	000372-38-3	9
5		Xenon	132	Xe	007440-63-3	9



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF120815\
Data File : BF083624.D
Acq On : 9 Dec 2015 6:08
Operator : UM/IZ
Sample : G4647-05
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
BRIDGE14-3(0-2)

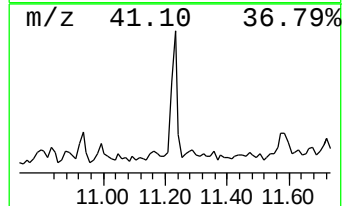
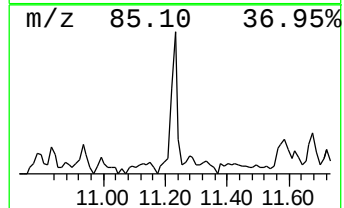
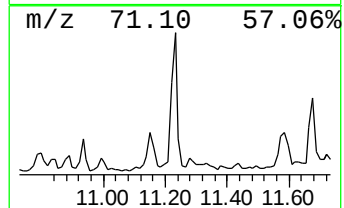
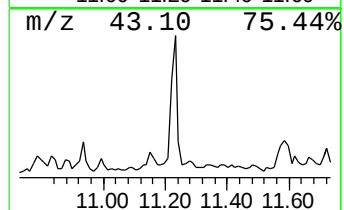
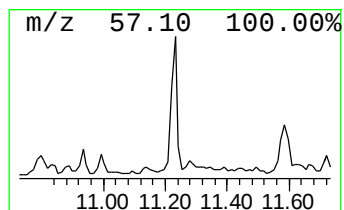
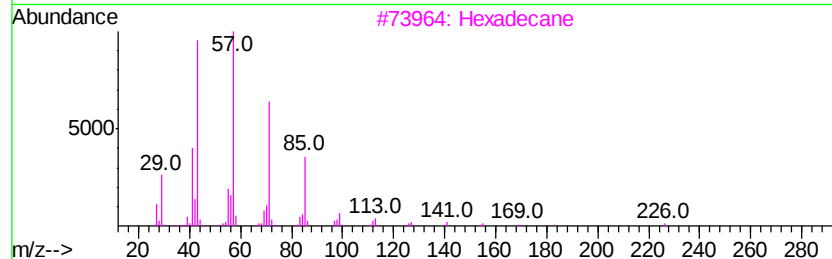
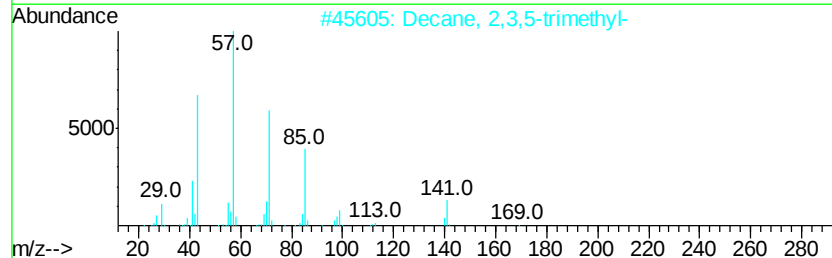
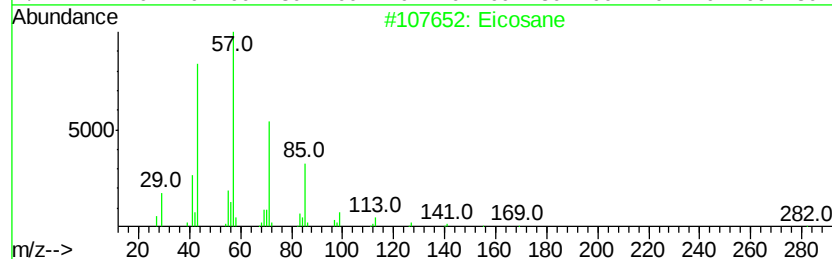
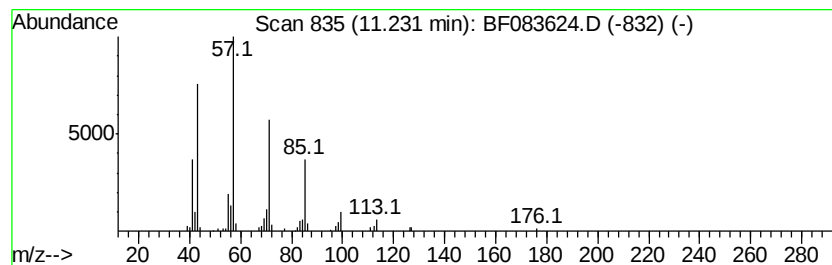
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF120815.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Eicosane Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.23	2.11 ng	124190	Acenaphthene-d10	10.38

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane		282	C20H42	000112-95-8	91
2	Decane, 2,3,5-trimethyl-		184	C13H28	062238-11-3	90
3	Hexadecane		226	C16H34	000544-76-3	90
4	Tetradecane		198	C14H30	000629-59-4	86
5	Pentadecane		212	C15H32	000629-62-9	86



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF120815\
Data File : BF083624.D
Acq On : 9 Dec 2015 6:08
Operator : UM/IZ
Sample : G4647-05
Misc :
ALS Vial : 32 Sample Multiplier: 1

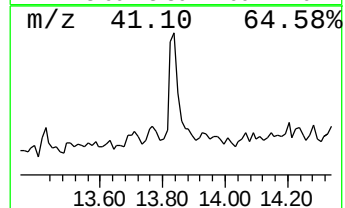
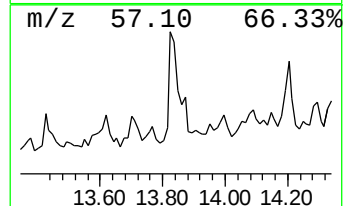
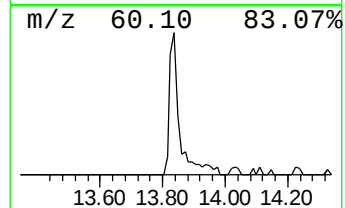
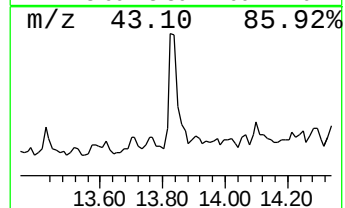
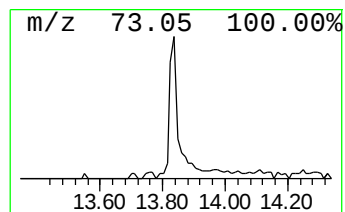
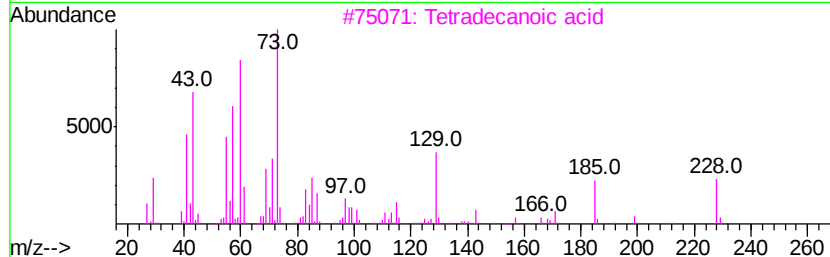
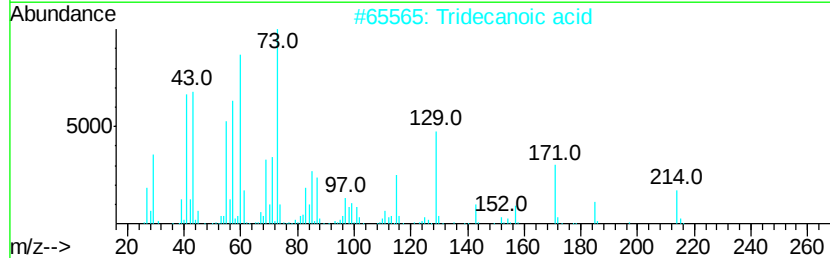
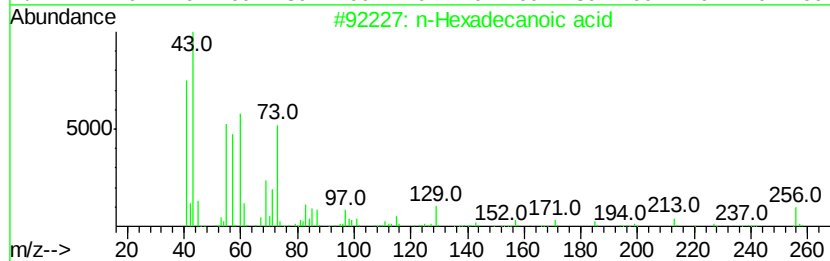
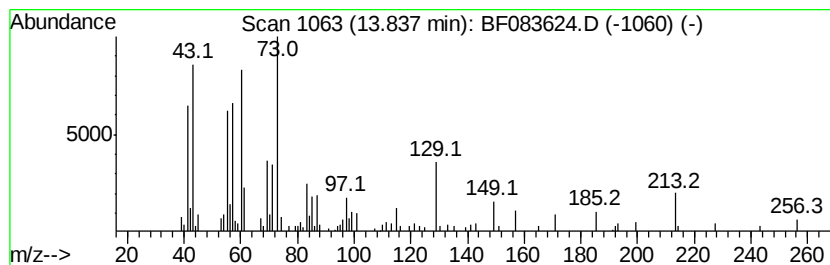
Instrument :
BNA_F
ClientSampleId :
BRIDGE14-3(0-2)

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 n-Hexadecanoic acid Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD		R.T.		
13.84	4.09 ng	183410	Phenanthrene-d10		12.79		
Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	96
2			Tridecanoic acid	214	C13H26O2	000638-53-9	76
3			Tetradecanoic acid	228	C14H28O2	000544-63-8	72
4			n-Decanoic acid	172	C10H20O2	000334-48-5	53
5			Undecanoic acid	186	C11H22O2	000112-37-8	53



Data Path : Z:\HPCHEM1\BNA F\DATA\BF120815\
Data File : BF083624.D
Acq On : 9 Dec 2015 6:08
Operator : UM/IZ
Sample : G4647-05
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
BRIDGE14-3(0-2)

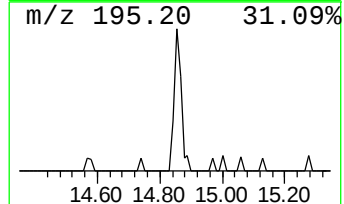
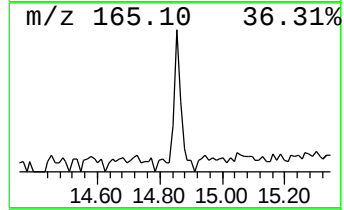
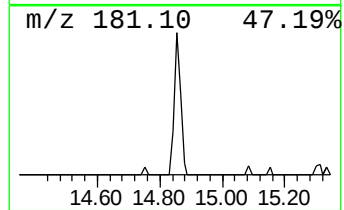
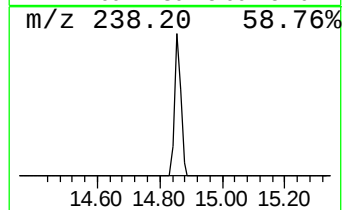
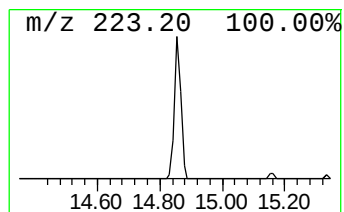
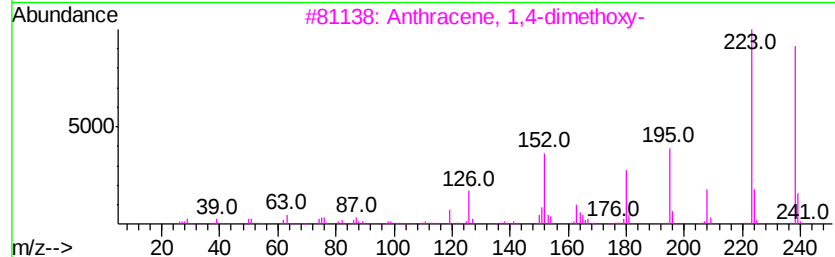
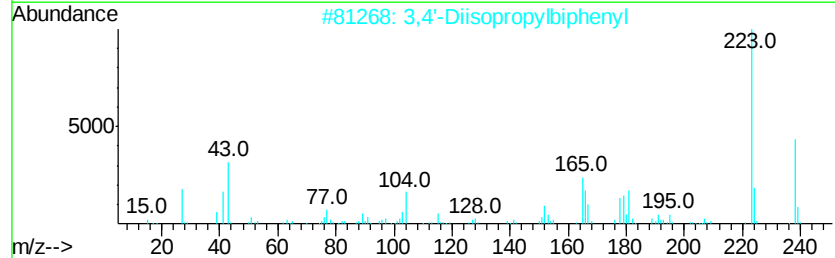
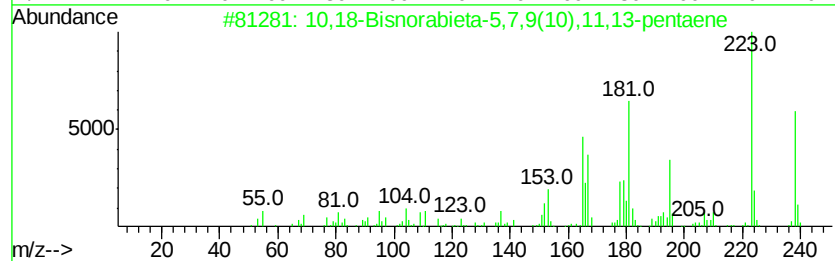
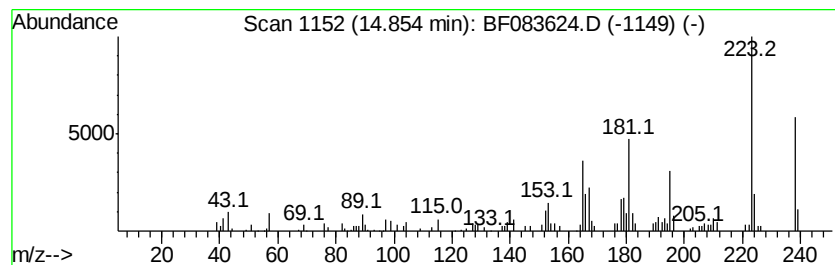
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF120815.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 10,18-Bisnorabieta-5,7,9(10... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.85	2.53 ng	113329	Phenanthrene-d10	12.79

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			10,18-Bisnorabieta-5,7,9(10),11,...	238	C18H22	006566-19-4	99
2			3,4'-Diisopropylbiphenyl	238	C18H22	061434-46-6	89
3			Anthracene, 1,4-dimethoxy-	238	C16H14O2	013076-29-4	58
4			2-Propenoic acid, 3-(3,4,5-trime...	238	C12H14O5	000090-50-6	53
5			4,4'-Diisopropylbiphenyl	238	C18H22	018970-30-4	52



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF120815\
Data File : BF083624.D
Acq On : 9 Dec 2015 6:08
Operator : UM/IZ
Sample : G4647-05
Misc :
ALS Vial : 32 Sample Multiplier: 1

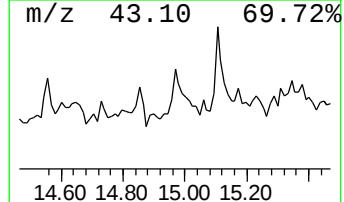
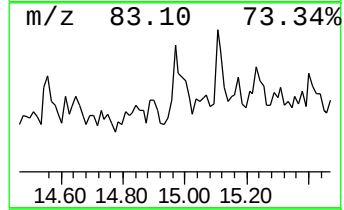
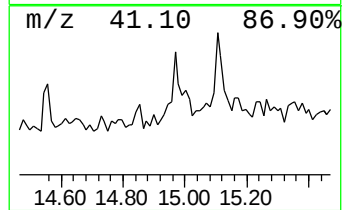
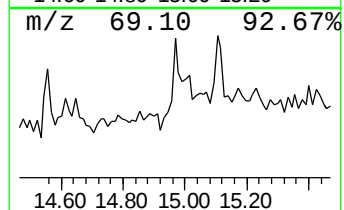
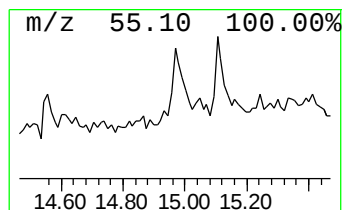
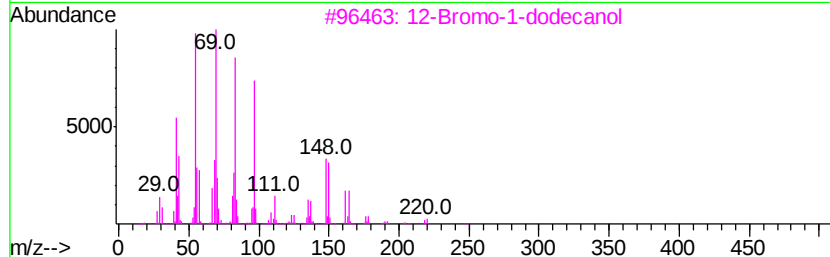
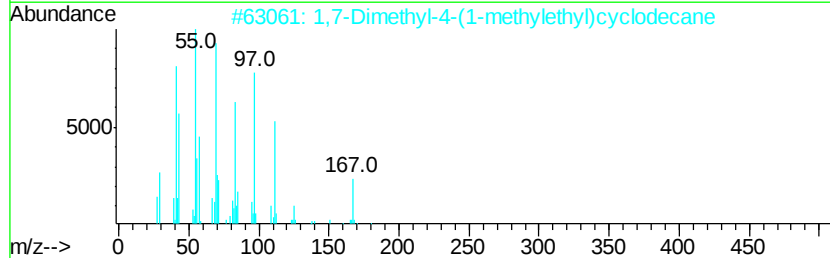
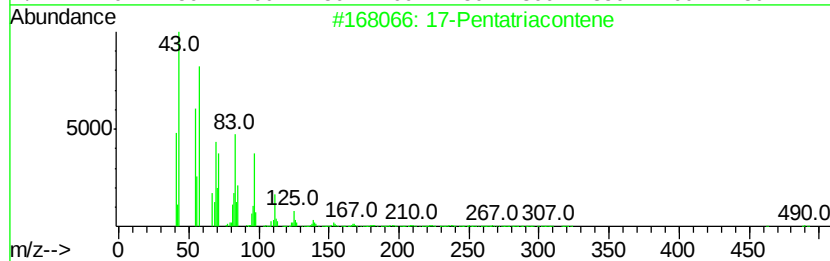
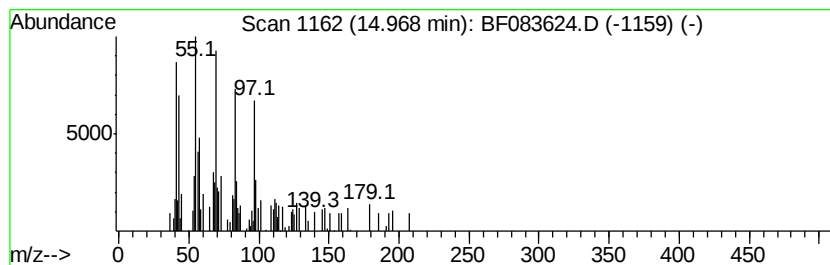
Instrument :
BNA_F
ClientSampleId :
BRIDGE14-3(0-2)

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 17-Pentatriacontene Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.97	2.22 ng	107924	Chrysene-d12	16.99
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	17-Pentatriacontene	491 C35H70	006971-40-0	72
2	1,7-Dimethyl-4-(1-methylethyl)cyclo...	210 C15H30	000645-10-3	53
3	12-Bromo-1-dodecanol	264 C12H25BrO	003344-77-2	50
4	2-Hexenal, (E)-	98 C6H10O	006728-26-3	49
5	3-Hexene, 2-methyl-, (Z)-	98 C7H14	015840-60-5	38



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF120815\
Data File : BF083624.D
Acq On : 9 Dec 2015 6:08
Operator : UM/IZ
Sample : G4647-05
Misc :
ALS Vial : 32 Sample Multiplier: 1

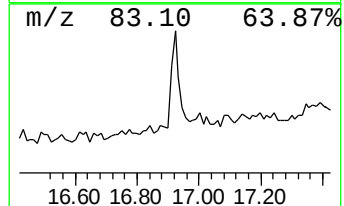
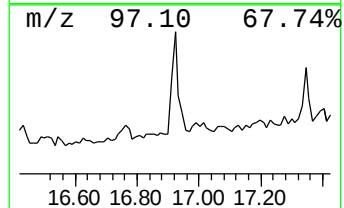
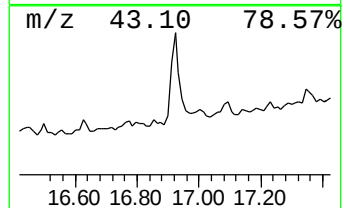
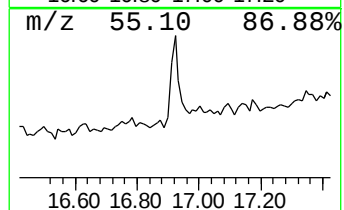
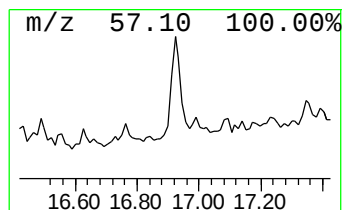
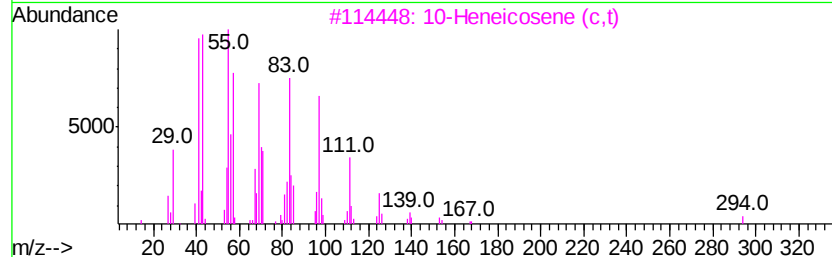
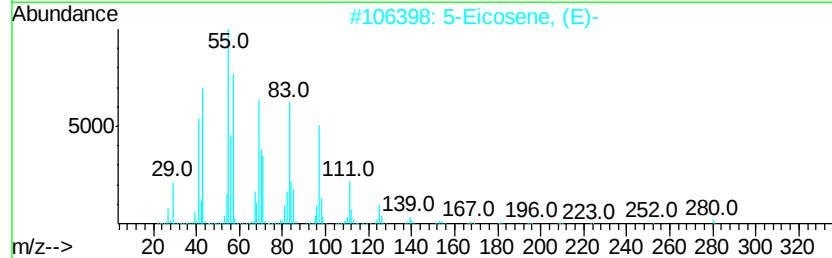
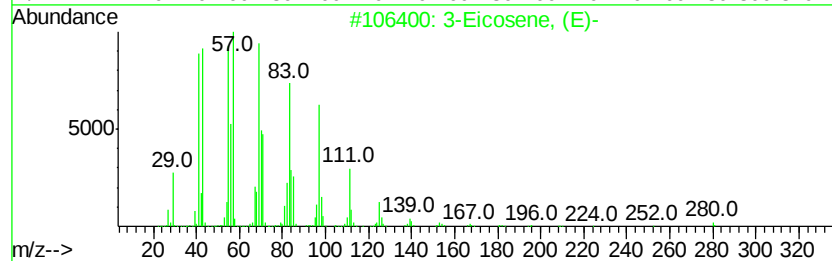
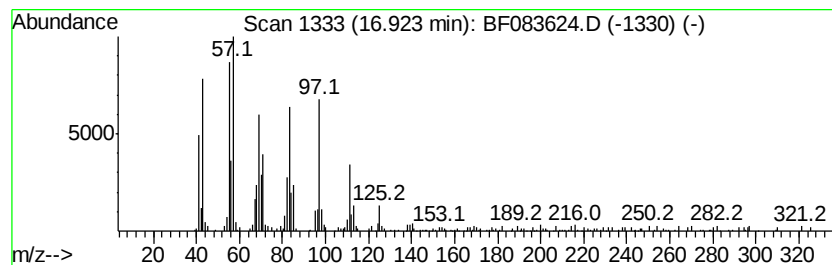
Instrument :
BNA_F
ClientSampleId :
BRIDGE14-3(0-2)

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 3-Eicosene, (E)- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD		R.T.
16.92	6.57 ng	319369	Chrysene-d12		16.99
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Eicosene, (E)-	280	C20H40	074685-33-9	97
2	5-Eicosene, (E)-	280	C20H40	074685-30-6	93
3	10-Heneicosene (c,t)	294	C21H42	095008-11-0	91
4	9-Eicosene, (E)-	280	C20H40	074685-29-3	91
5	1-Heneicosyl formate	340	C22H44O2	077899-03-7	91



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF120815\
Data File : BF083624.D
Acq On : 9 Dec 2015 6:08
Operator : UM/IZ
Sample : G4647-05
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
BRIDGE14-3(0-2)

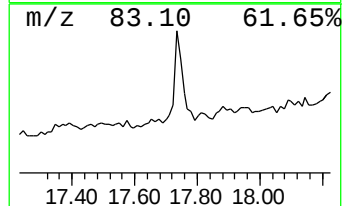
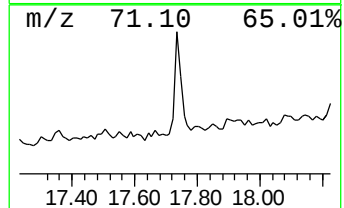
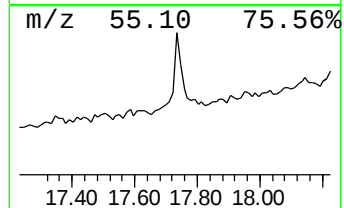
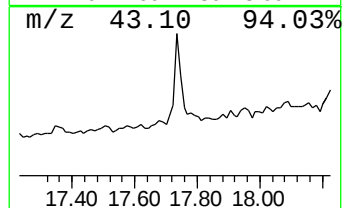
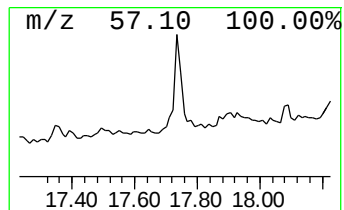
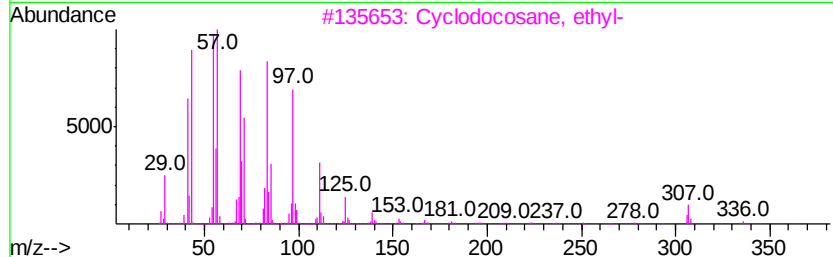
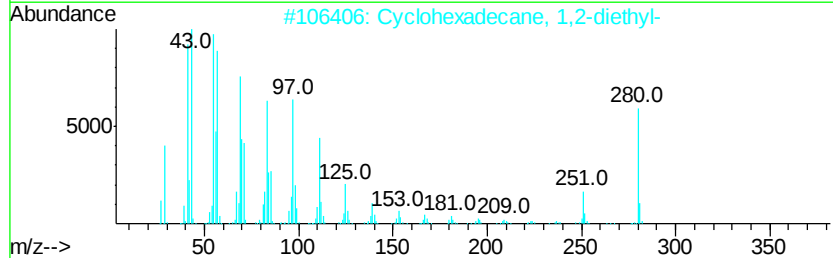
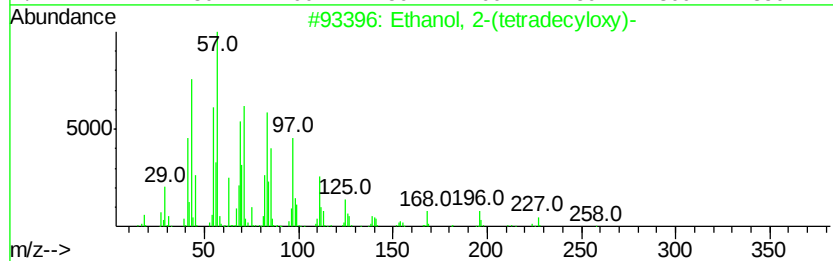
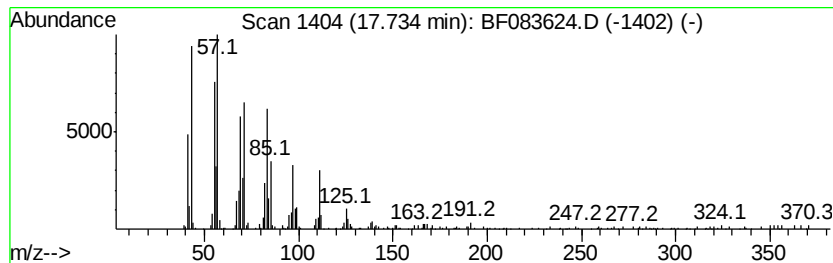
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF120815.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Ethanol, 2-(tetradecyloxy)- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.73	6.11 ng	296565	Chrysene-d12	16.99

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	91
2			Cyclohexadecane, 1,2-diethyl-	280	C20H40	1000155-85-3	90
3			Cyclodocosane, ethyl-	336	C24H48	1000151-22-6	89
4			1-Octadecanol	270	C18H38O	000112-92-5	76
5			17-Pentatriacontene	491	C35H70	006971-40-0	74



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF120815\
Data File : BF083624.D
Acq On : 9 Dec 2015 6:08
Operator : UM/IZ
Sample : G4647-05
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
BRIDGE14-3(0-2)

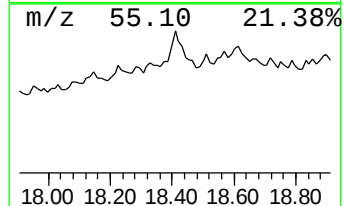
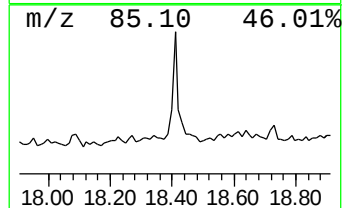
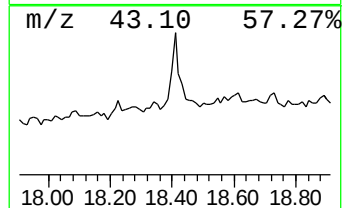
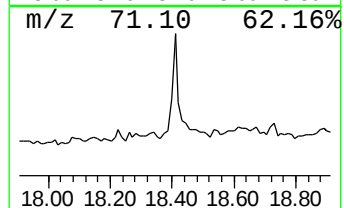
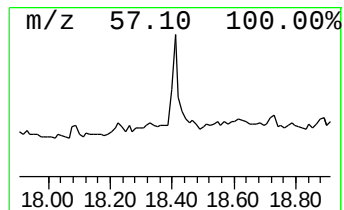
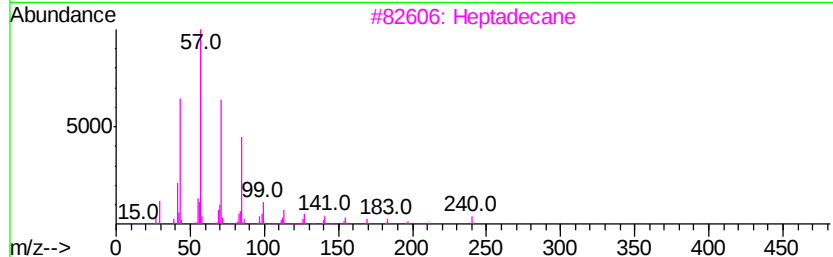
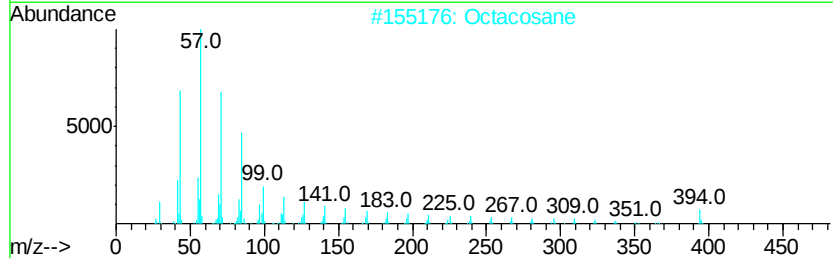
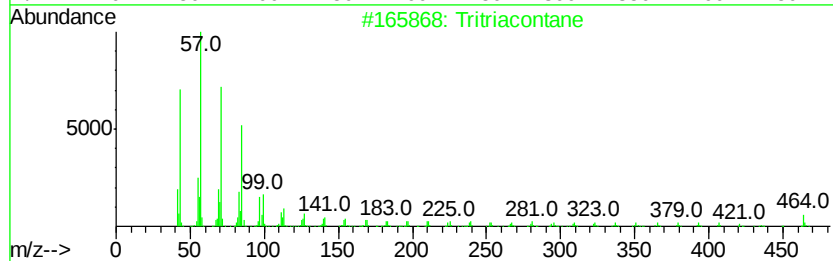
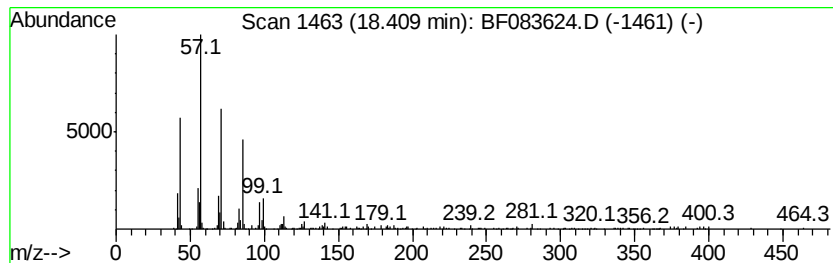
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF120815.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Tritriacontane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.41	6.56 ng	207641	Perylene-d12	18.64

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tritriacontane	465	C33H68	000630-05-7	90
2			Octacosane	394	C28H58	000630-02-4	90
3			Heptadecane	240	C17H36	000629-78-7	83
4			Pentadecane	212	C15H32	000629-62-9	81
5			Eicosane	282	C20H42	000112-95-8	80



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF120815\
Data File : BF083624.D
Acq On : 9 Dec 2015 6:08
Operator : UM/IZ
Sample : G4647-05
Misc :
ALS Vial : 32 Sample Multiplier: 1

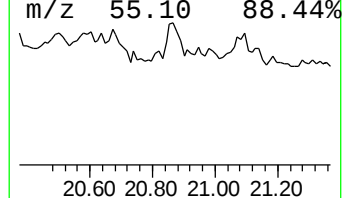
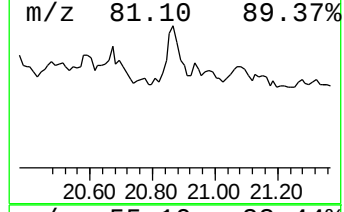
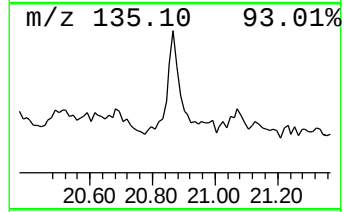
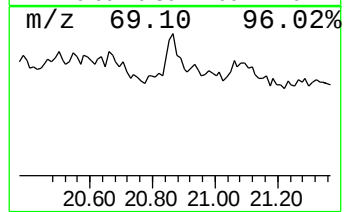
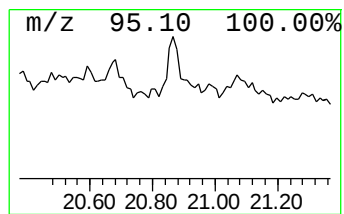
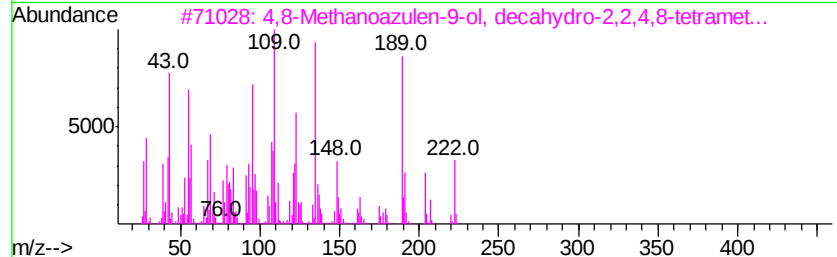
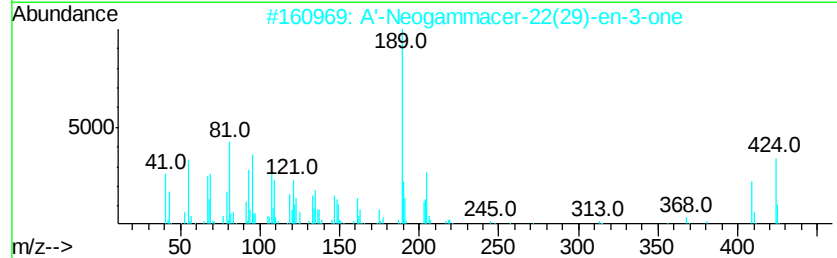
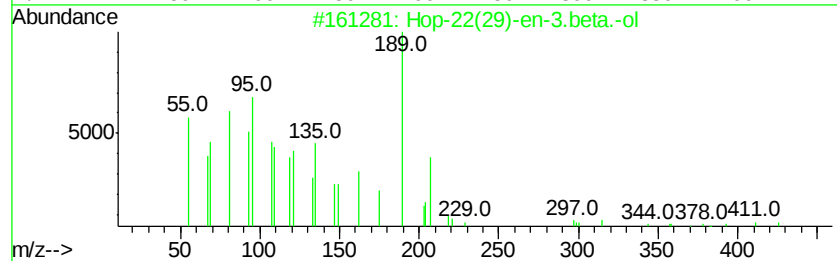
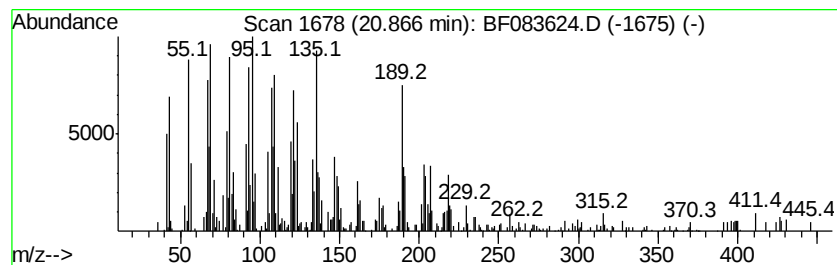
Instrument :
BNA_F
ClientSampleId :
BRIDGE14-3(0-2)

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 Hop-22(29)-en-3.beta.-ol Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.87	11.75 ng	371943	Perylene-d12	18.64
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Hop-22(29)-en-3.beta.-ol	426 C30H500	058801-23-3 94
2		A'-Neogammacer-22(29)-en-3-one	424 C30H480	025615-11-6 93
3		4,8-Methanoazulen-9-ol, decahvd...	222 C15H260	004586-22-5 58
4		1-Formyl-2,2,6-trimethyl-3-cis-(...	220 C15H240	1000144-10-1 55
5		Cedran-diol, 8S,14-	238 C15H2602	062600-05-9 55



Data Path : Z:\HPCHEM1\BNA F\DATA\BF120815\
Data File : BF083624.D
Acq On : 9 Dec 2015 6:08
Operator : UM/IZ
Sample : G4647-05
Misc :
ALS Vial : 32 Sample Multiplier: 1

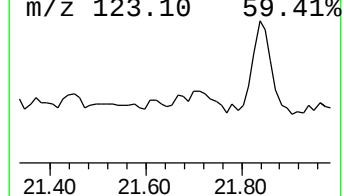
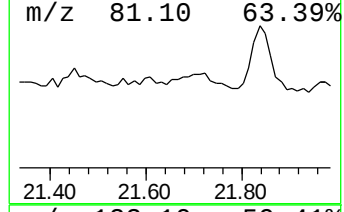
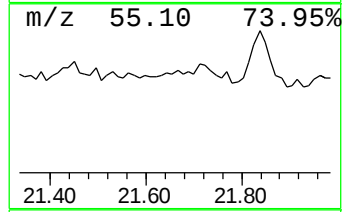
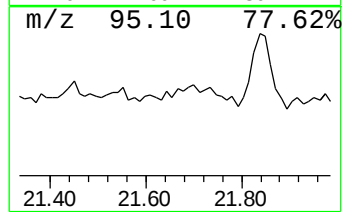
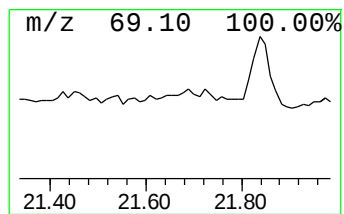
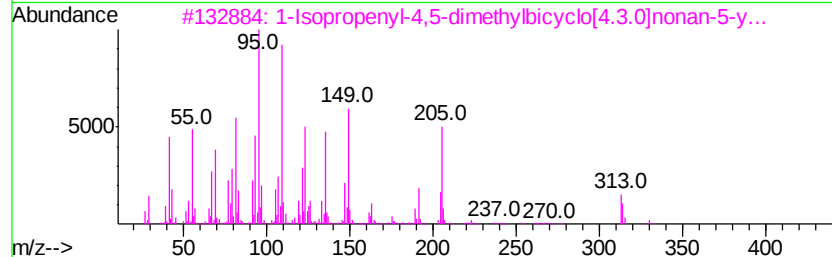
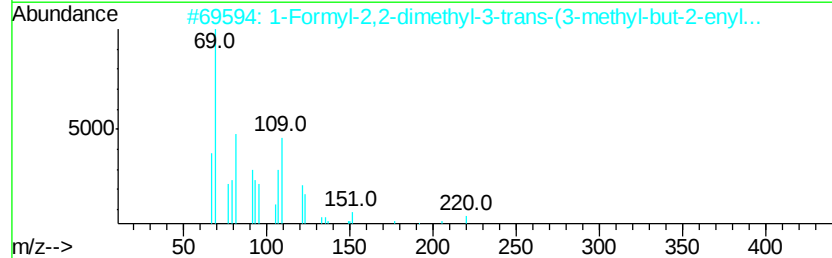
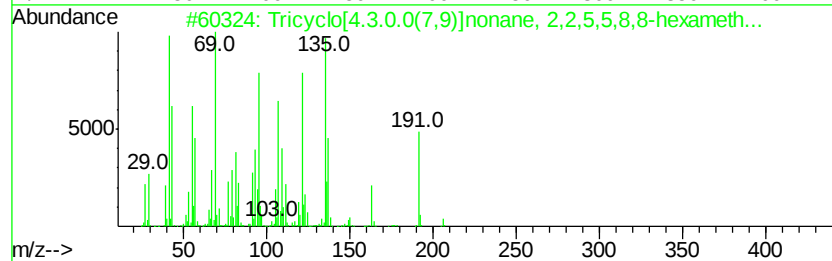
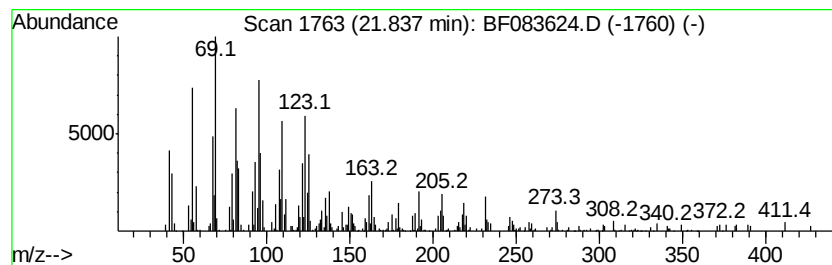
Instrument :
BNA_F
ClientSampleId :
BRIDGE14-3(0-2)

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 unknown21.84 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.84	17.80 ng	563539	Perylene-d12	18.64
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Tricyclo[4.3.0.0(7,9)]nonane, 2,...	206 C15H26	054832-82-5 43
2		1-Formyl-2,2-dimethyl-3-trans-(3...	220 C15H24O	1000144-09-7 42
3		1-Isopropenyl-4,5-dimethylbicycl...	330 C21H30OS	1000195-85-4 30
4		1,6,10,14,18,22-Tetracosahexaen...	426 C30H50O	054159-46-5 25
5		6.beta.Bicyclo[4.3.0]nonane, 5.b...	332 C15H25I	1000195-85-9 20



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF120815\
Data File : BF083624.D
Acq On : 9 Dec 2015 6:08
Operator : UM/IZ
Sample : G4647-05
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
BRIDGE14-3(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
unknown1.85	1.85	10.3	ng	478613	1	5.70	933050	20.0
2-Pentanone, 4-hy...	3.53	14.6	ng	681693	1	5.70	933050	20.0
unknown5.40	5.40	63.4	ng	2957180	1	5.70	933050	20.0
Eicosane	11.23	2.1	ng	124190	3	10.38	1175520	20.0
n-Hexadecanoic acid	13.84	4.1	ng	183410	4	12.79	896461	20.0
10,18-Bisnorabiet...	14.85	2.5	ng	113329	4	12.79	896461	20.0
17-Pentatriacontene	14.97	2.2	ng	107924	5	16.99	971545	20.0
3-Eicosene, (E)-	16.92	6.6	ng	319369	5	16.99	971545	20.0
Ethanol, 2-(tetra...	17.73	6.1	ng	296565	5	16.99	971545	20.0
Tritriacontane	18.41	6.6	ng	207641	6	18.64	633247	20.0
Hop-22(29)-en-3.b...	20.87	11.8	ng	371943	6	18.64	633247	20.0
unknown21.84	21.84	17.8	ng	563539	6	18.64	633247	20.0