

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF011116\
 Data File : BF084381.D
 Acq On : 11 Jan 2016 15:54
 Operator : SJ/UM
 Sample : H1035-02
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 HAZ-COMP

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-BF123115.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	5.036	255	258	267	rVB	1115942	1484636	27.14%	2.524%
2	5.470	293	296	298	rBV	5205456	5363210	98.05%	9.119%
3	6.499	383	386	389	rBV	4836129	4760296	87.02%	8.094%
4	6.625	394	397	400	rBV	5484264	4846349	88.60%	8.240%
5	6.853	414	417	419	rBV	1985254	1614916	29.52%	2.746%
6	7.013	428	431	433	rVB	4169741	4193470	76.66%	7.130%
7	7.425	464	467	469	rBV	2635828	3575600	65.37%	6.079%
8	7.539	475	477	482	rBV	48368	80676	1.47%	0.137%
9	8.145	527	530	532	rBV	1730119	2013050	36.80%	3.423%
10	9.219	621	624	626	rBV	6281942	5470128	100.00%	9.301%
11	9.608	656	658	661	rBV	620980	821972	15.03%	1.398%
12	9.825	673	677	681	rBV2	80047	161437	2.95%	0.274%
13	9.893	681	683	685	rBV	2496785	2216345	40.52%	3.768%
14	10.019	691	694	696	rBV	286966	284920	5.21%	0.484%
15	10.328	720	721	723	rVB	192602	147909	2.70%	0.251%
16	10.545	737	740	744	rBV	62983	121096	2.21%	0.206%
17	10.682	749	752	754	rVV	3278245	3147631	57.54%	5.352%
18	10.716	754	755	761	rVB2	84861	127829	2.34%	0.217%
19	10.796	761	762	767	rBV	134696	240500	4.40%	0.409%
20	11.254	799	802	803	rBV	118108	116508	2.13%	0.198%
21	11.379	810	813	815	rBV	2254435	2007004	36.69%	3.412%
22	11.436	816	818	822	rVB2	59457	87074	1.59%	0.148%
23	11.608	830	833	837	rBV3	34550	76395	1.40%	0.130%
24	11.676	837	839	840	rVV	71994	77597	1.42%	0.132%
25	11.699	840	841	843	rVB	241464	180191	3.29%	0.306%
26	11.882	855	857	858	rVB2	111068	122463	2.24%	0.208%
27	12.111	875	877	880	rVB2	252332	308948	5.65%	0.525%
28	12.282	887	892	897	rVB6	37878	86507	1.58%	0.147%
29	12.351	897	898	901	rBV3	84180	77834	1.42%	0.132%
30	12.408	901	903	904	rBV	178187	165246	3.02%	0.281%
31	12.442	904	906	907	rVV2	65336	106728	1.95%	0.181%
32	12.488	907	910	914	rVV4	119233	245256	4.48%	0.417%
33	12.591	914	919	920	rVV	101924	287711	5.26%	0.489%
34	12.614	920	921	923	rVB2	148398	132344	2.42%	0.225%

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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

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Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	12.682	923	927	929	rBV2	738267	1027356	18.78%	1.747%
36	12.785	934	936	940	rBV3	790881	1413292	25.84%	2.403%
37	12.854	940	942	945	rVV	237603	360293	6.59%	0.613%
38	12.899	945	946	949	rVV	122246	140322	2.57%	0.239%
39	12.968	949	952	954	rVB	3957071	4047339	73.99%	6.882%
40	13.197	969	972	973	rBV2	31305	63411	1.16%	0.108%
41	13.288	978	980	983	rVV2	130562	131447	2.40%	0.223%
42	13.448	990	994	999	rBV	1279314	1332503	24.36%	2.266%
43	13.620	1007	1009	1011	rBV2	48405	63832	1.17%	0.109%
44	13.860	1028	1030	1036	rBV	362031	693425	12.68%	1.179%
45	14.008	1041	1043	1046	rBV2	1116360	1326673	24.25%	2.256%
46	14.500	1083	1086	1095	rBV2	156110	370030	6.76%	0.629%
47	14.865	1116	1118	1120	rVB	51867	61883	1.13%	0.105%
48	15.025	1130	1132	1137	rVB2	42216	91379	1.67%	0.155%
49	15.117	1137	1140	1141	rBV	79184	96925	1.77%	0.165%
50	15.437	1166	1168	1171	rBV	826422	1165406	21.30%	1.981%
51	15.677	1186	1189	1191	rBV3	32116	61465	1.12%	0.105%
52	15.894	1205	1208	1211	rVV	126412	183742	3.36%	0.312%
53	15.963	1211	1214	1217	rVV4	27985	80150	1.47%	0.136%
54	16.031	1217	1220	1223	rVB3	135897	247323	4.52%	0.421%
55	16.546	1262	1265	1272	rVB	39798	76605	1.40%	0.130%
56	16.843	1288	1291	1297	rVB2	29871	75294	1.38%	0.128%
57	16.946	1297	1300	1303	rBV	36470	83513	1.53%	0.142%
58	17.151	1313	1318	1324	rVB3	90275	215700	3.94%	0.367%
59	17.414	1337	1341	1343	rBV5	41783	123735	2.26%	0.210%
60	17.471	1343	1346	1350	rVB5	38278	108251	1.98%	0.184%
61	17.608	1355	1358	1362	rVB4	36627	97967	1.79%	0.167%
62	17.780	1369	1373	1379	rBV5	43001	113420	2.07%	0.193%
63	18.374	1422	1425	1431	rVB8	35391	118875	2.17%	0.202%
64	19.369	1508	1512	1517	rVB7	28152	93087	1.70%	0.158%

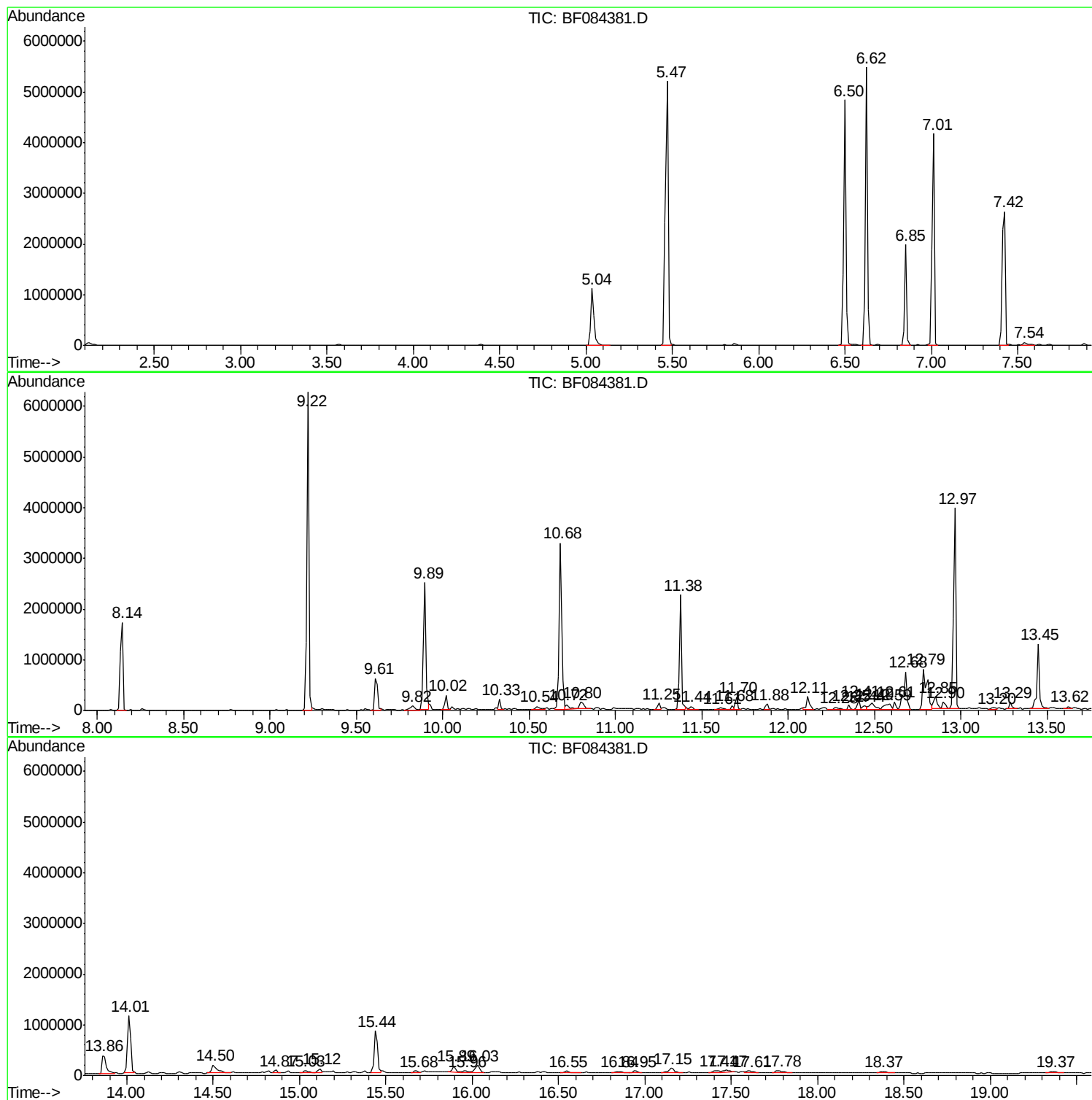
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ClientSampled :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF123115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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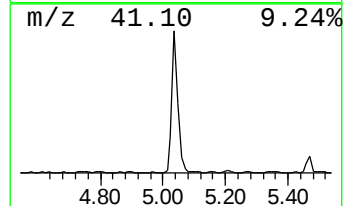
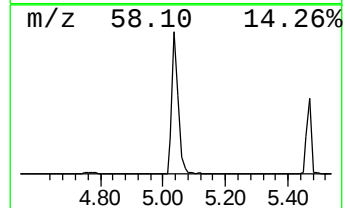
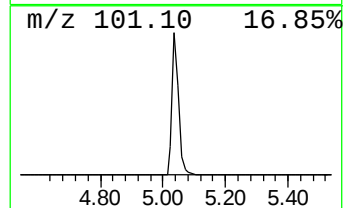
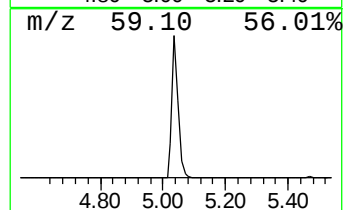
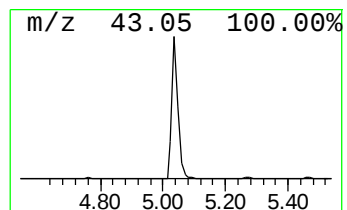
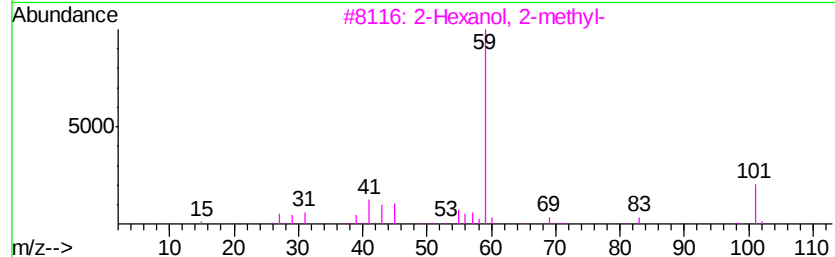
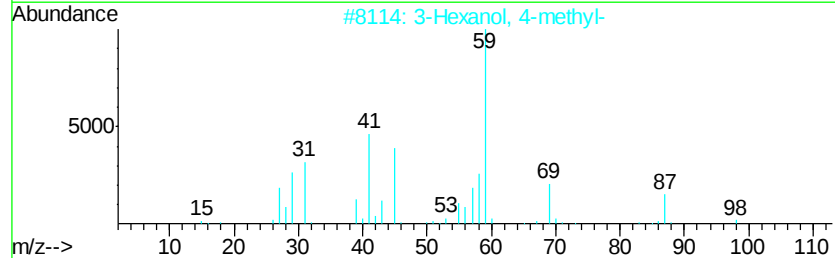
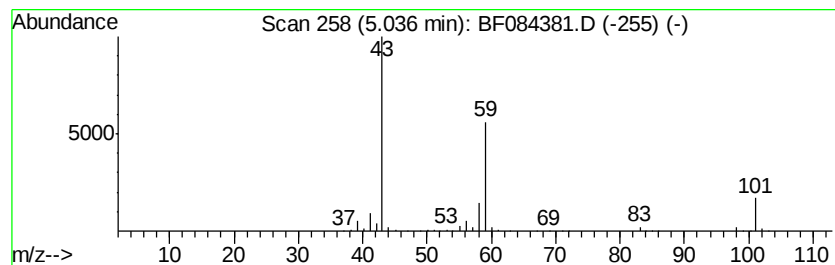
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.04	18.39 ng	1484640	1,4-Dichlorobenzene-d4	6.85

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50	
2	3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33	
3	2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	28	
4	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	25	
5	Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	25	



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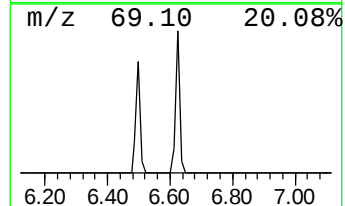
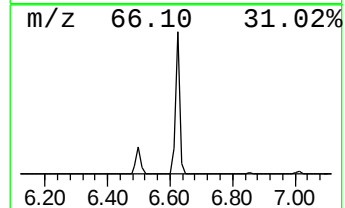
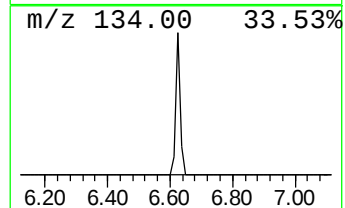
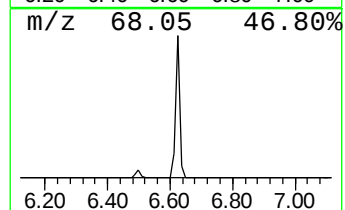
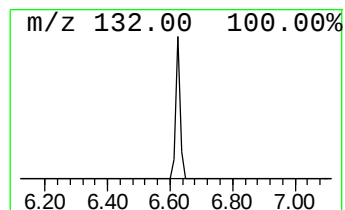
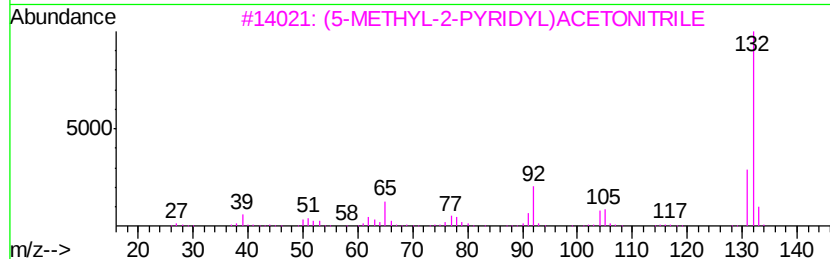
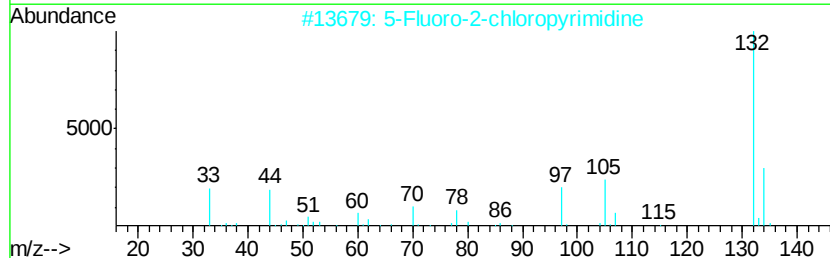
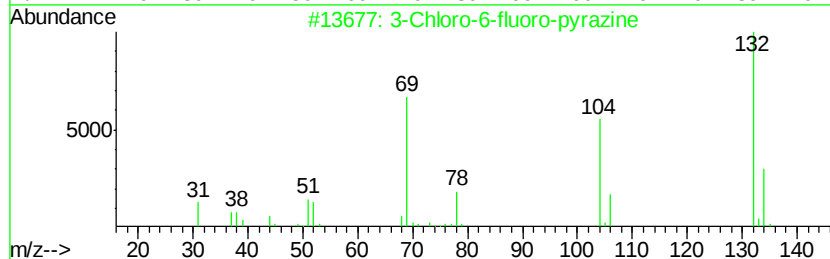
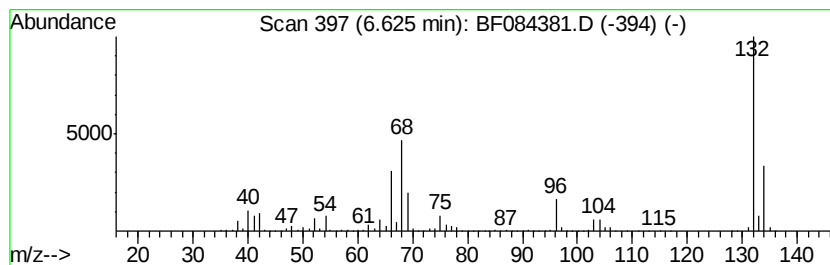
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown6.62 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.62	60.02 ng	4846350	1,4-Dichlorobenzene-d4	6.85

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
3		(5-METHYL-2-PYRIDYL)ACETONITRILE	132	C8H8N2	1000241-93-9	10
4		1,2,3-Trifluorobenzene	132	C6H3F3	001489-53-8	9
5		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9



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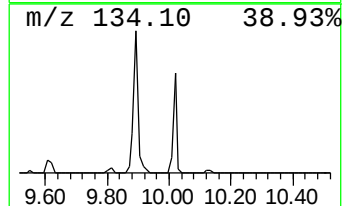
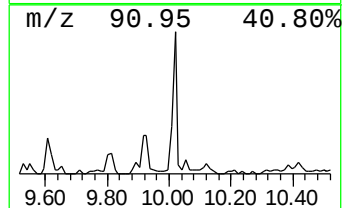
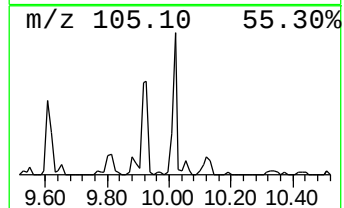
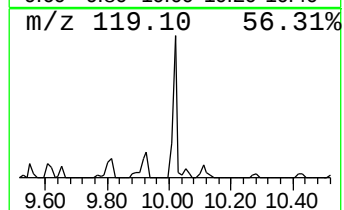
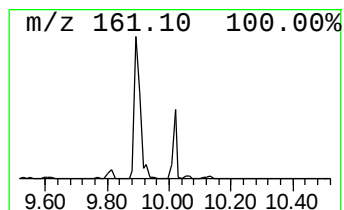
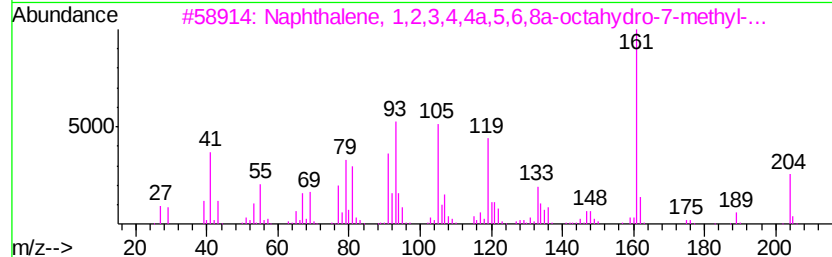
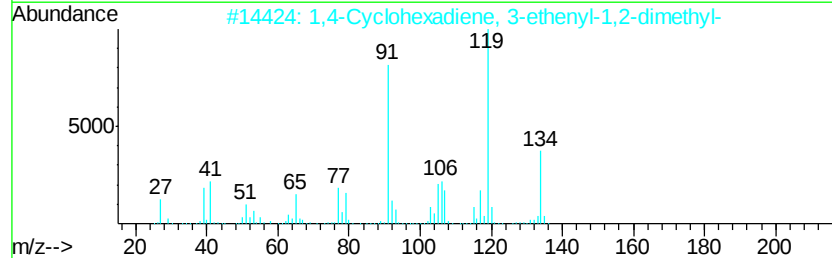
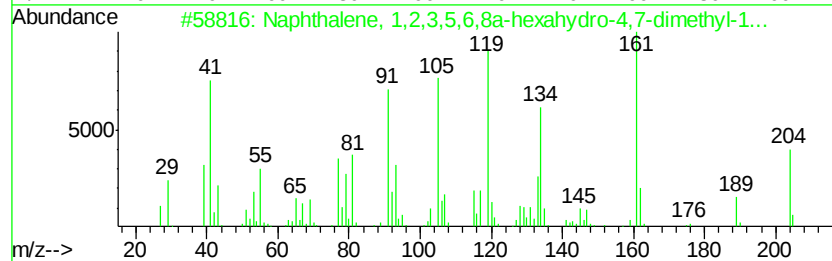
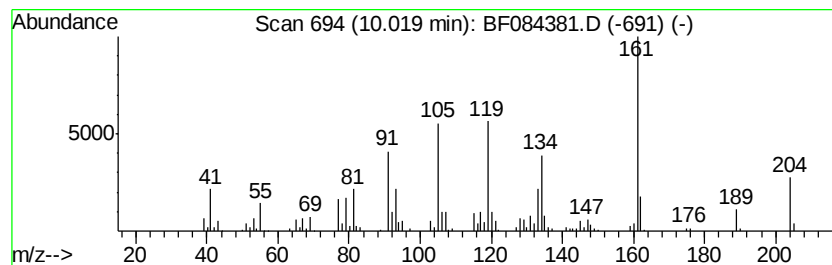
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Naphthalene, 1,2,3,5,6,8a-h... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.02	2.57 ng	284920	Acenaphthene-d10	9.89

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,2,3,5,6,8a-hexahy...	204	C15H24	000483-76-1	98
2		1,4-Cyclohexadiene, 3-ethenyl-1,...	134	C10H14	062338-57-2	83
3		Naphthalene, 1,2,3,4,4a,5,6,8a-o...	204	C15H24	030021-74-0	81
4		Naphthalene, 1,2,4a,5,8,8a-hexah...	204	C15H24	000523-47-7	76
5		Copaene	204	C15H24	003856-25-5	62



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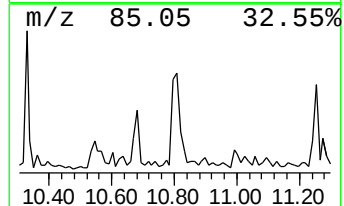
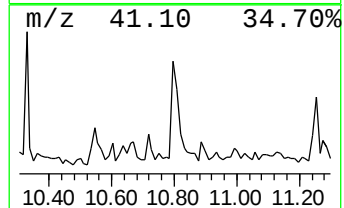
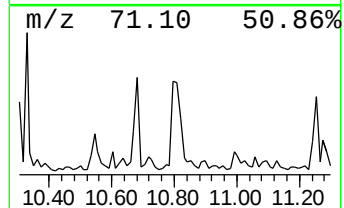
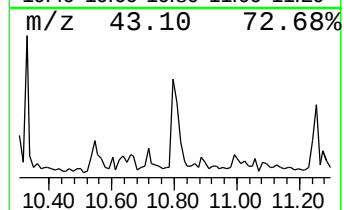
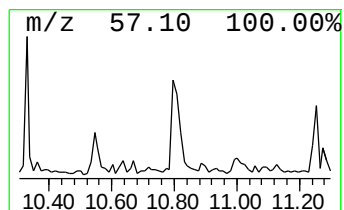
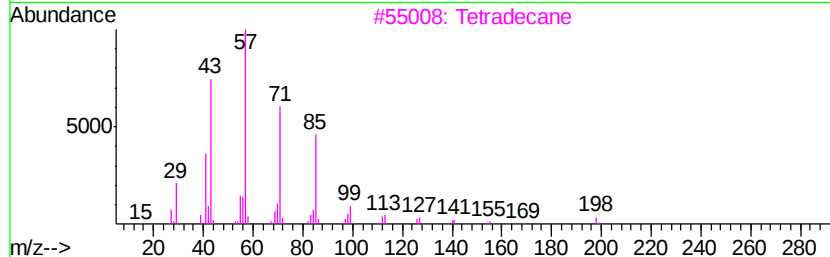
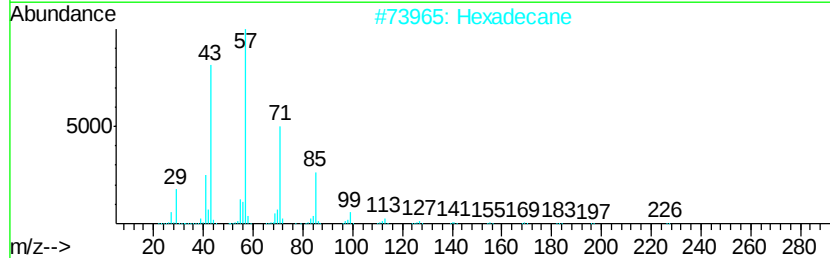
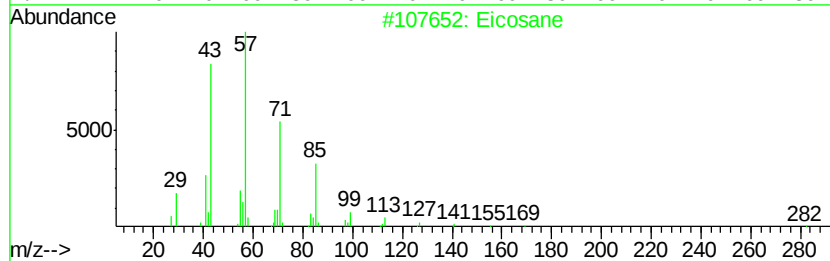
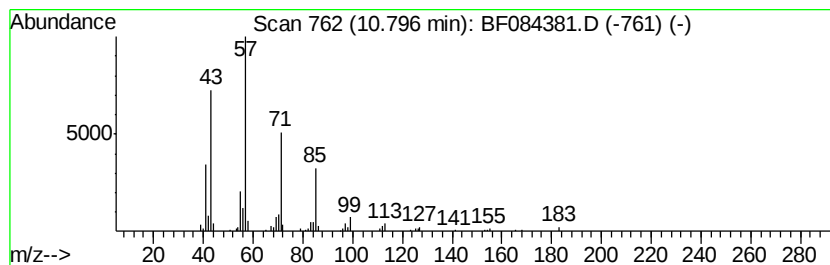
Instrument :
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ClientSampleId :
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Eicosane Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD		R.T.
10.80	2.40 ng	240500	Phenanthrene-d10		11.38
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane	282	C20H42	000112-95-8	91
2	Hexadecane	226	C16H34	000544-76-3	90
3	Tetradecane	198	C14H30	000629-59-4	86
4	Tridecane, 7-propyl-	226	C16H34	055045-09-5	80
5	Dodecane	170	C12H26	000112-40-3	78



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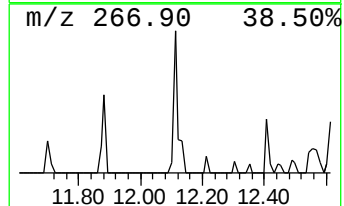
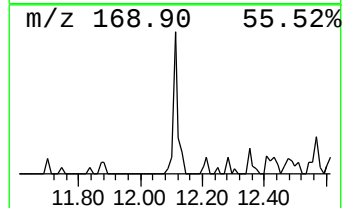
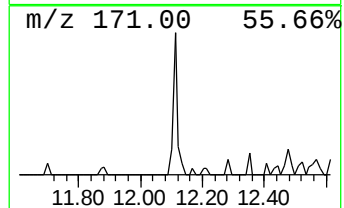
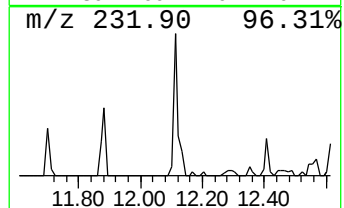
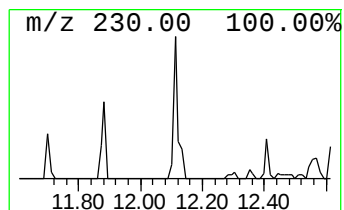
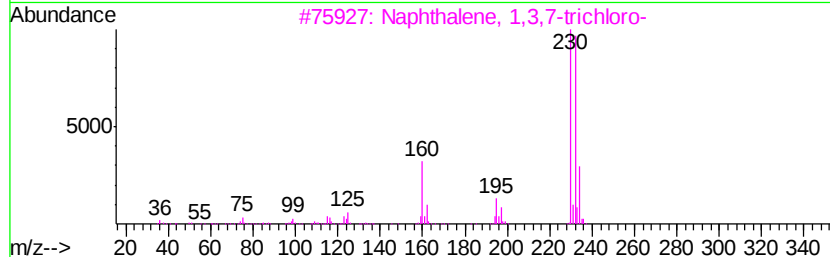
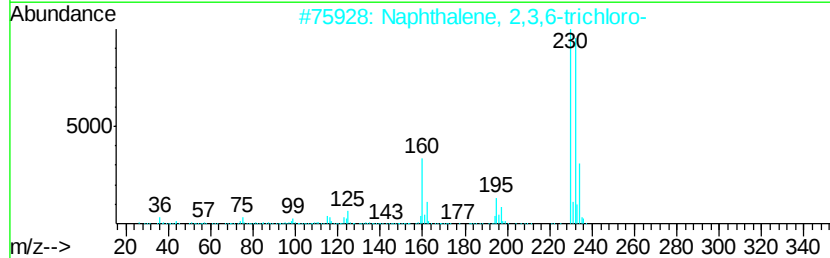
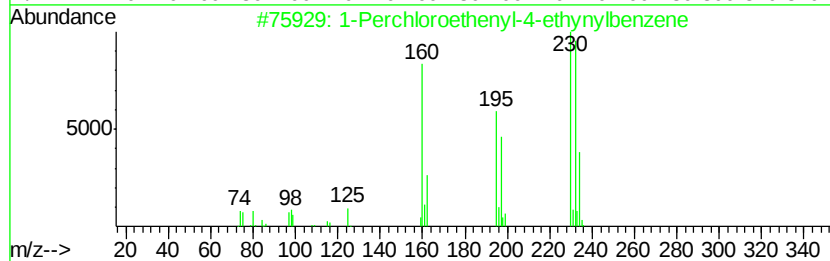
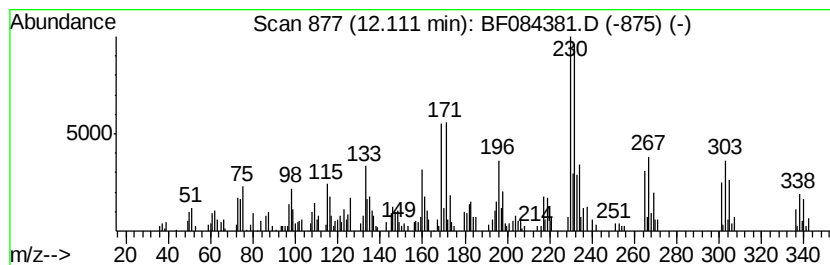
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 1-Perchloroethenyl-4-ethyny... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.11	3.08 ng	308948	Phenanthrene-d10	11.38

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Perchloroethenyl-4-ethynylbenzene	230	C10H5Cl3	1000283-76-9	91
2		Naphthalene, 2,3,6-trichloro-	230	C10H5Cl3	055720-40-6	55
3		Naphthalene, 1,3,7-trichloro-	230	C10H5Cl3	055720-37-1	55
4		Phenol, 2,3,4,5-tetrachloro-	230	C6H2Cl4O	004901-51-3	45
5		1,6-Methano-1H-indene, 2,3,3a,4,...	336	C10H6Cl6	056641-38-4	42



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF011116\
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Acq On : 11 Jan 2016 15:54
Operator : SJ/UM
Sample : H1035-02
Misc :
ALS Vial : 11 Sample Multiplier: 1

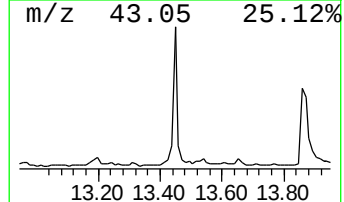
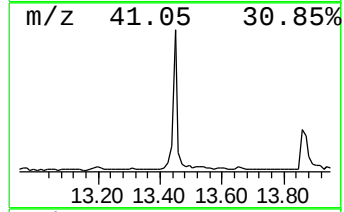
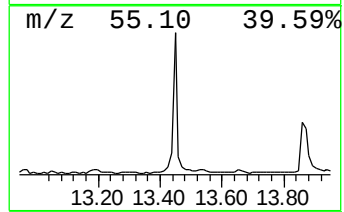
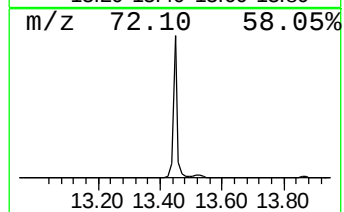
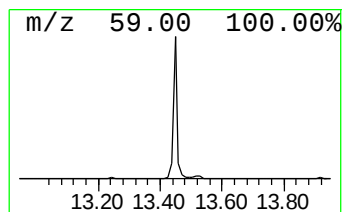
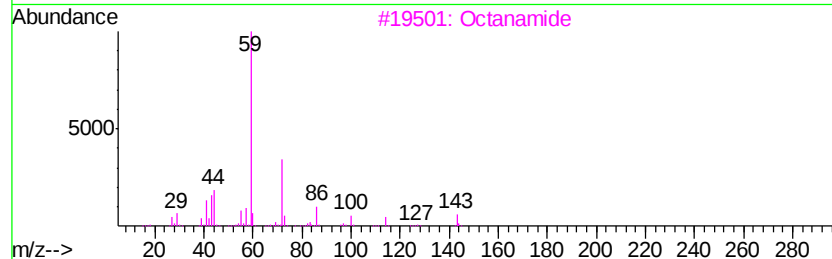
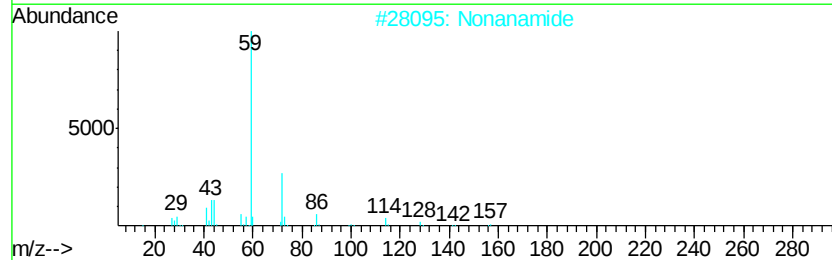
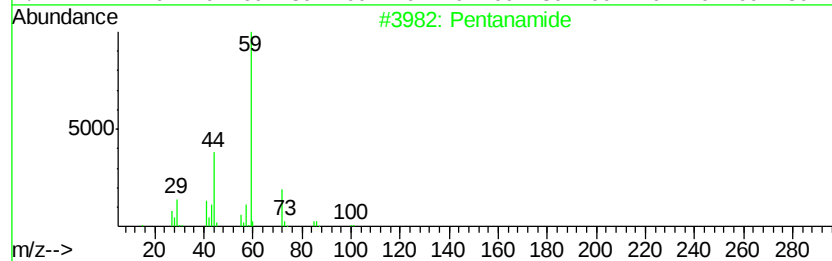
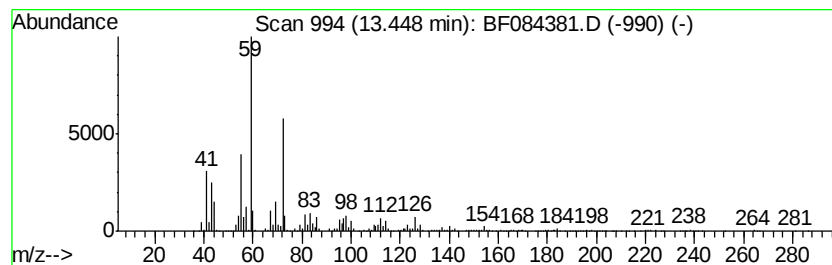
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Pentanamide Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.45	20.09 ng	1332500	Chrysene-d12	14.01
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Pentanamide	101	C5H11NO	000626-97-1 59
2	Nonanamide	157	C9H19NO	001120-07-6 53
3	Octanamide	143	C8H17NO	000629-01-6 53
4	Heptanamide, 4-ethyl-5-methyl-	171	C10H21NO	054789-40-1 50
5	1,7-Octanediol, 3,7-dimethyl-	174	C10H22O2	000107-74-4 43



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF011116\
Data File : BF084381.D
Acq On : 11 Jan 2016 15:54
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Sample : H1035-02
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Instrument :
BNA_F
ClientSampleId :
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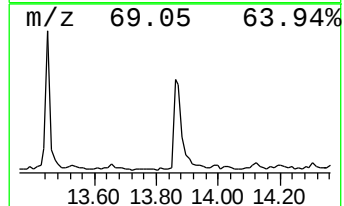
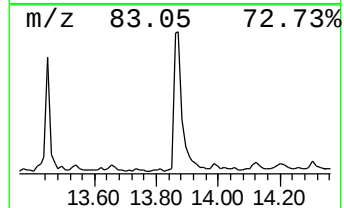
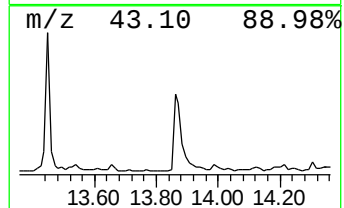
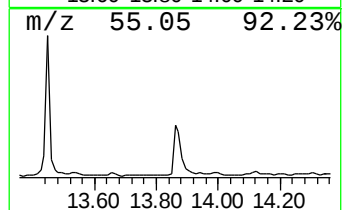
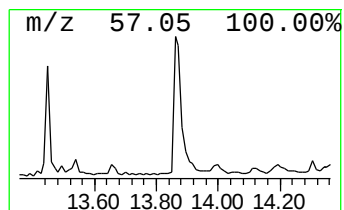
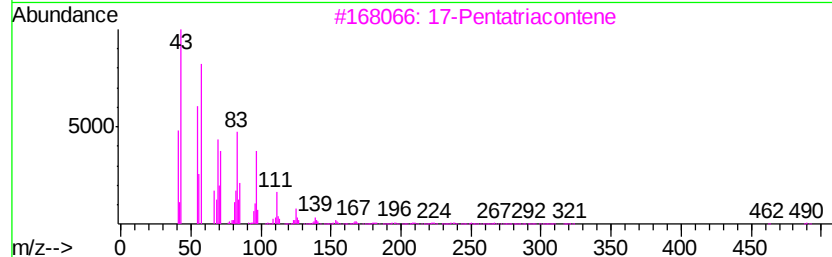
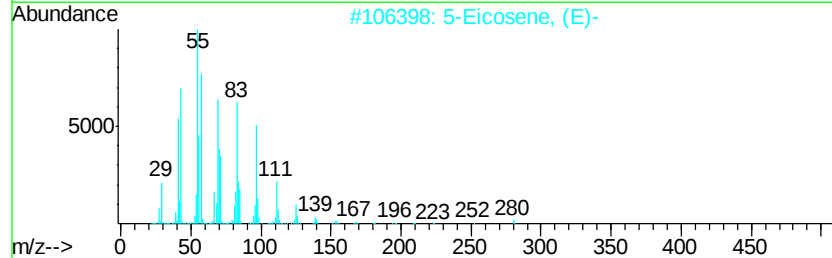
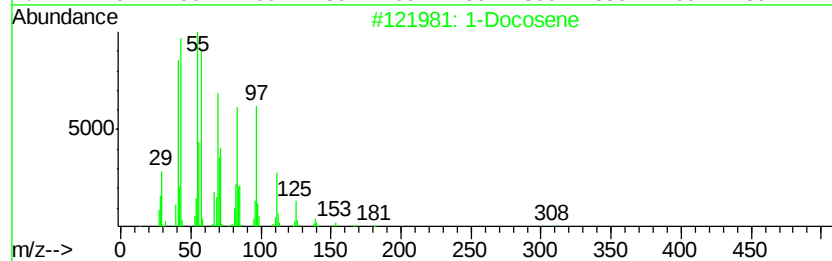
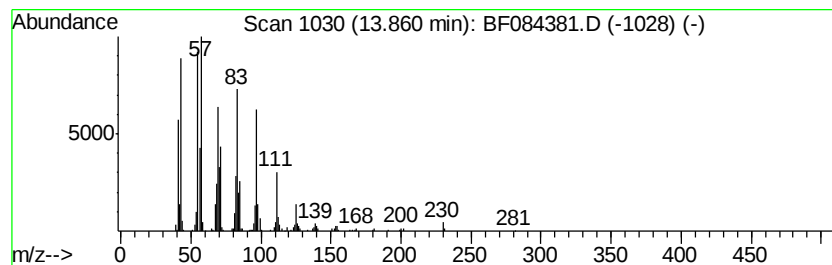
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 1-Docosene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.86	10.45 ng	693425	Chrysene-d12	14.01

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Docosene	308	C22H44	001599-67-3	94
2		5-Eicosene, (E)-	280	C20H40	074685-30-6	90
3		17-Pentatriacontene	491	C35H70	006971-40-0	90
4		3-Eicosene, (E)-	280	C20H40	074685-33-9	90
5		1-Nonadecene	266	C19H38	018435-45-5	90



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF011116\
Data File : BF084381.D
Acq On : 11 Jan 2016 15:54
Operator : SJ/UM
Sample : H1035-02
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
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ClientSampleId :
HAZ-COMP

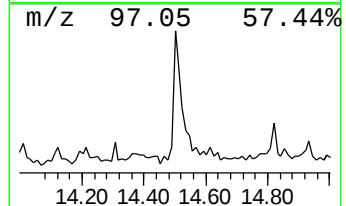
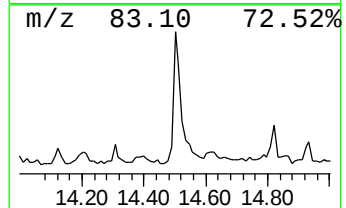
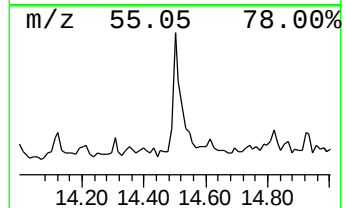
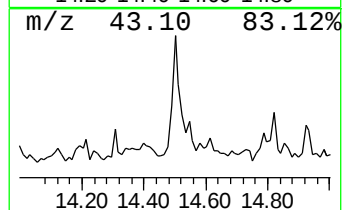
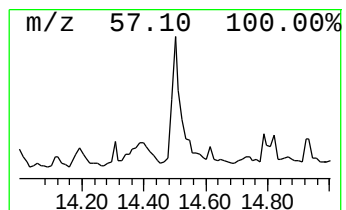
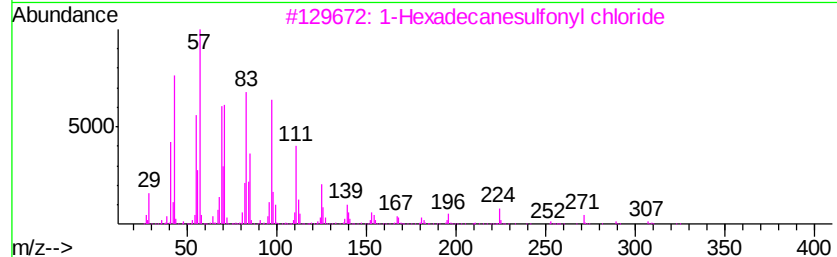
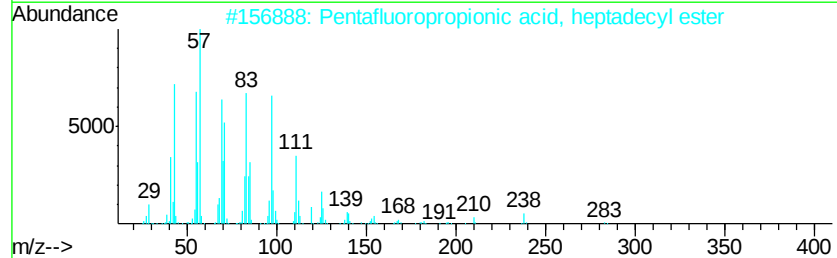
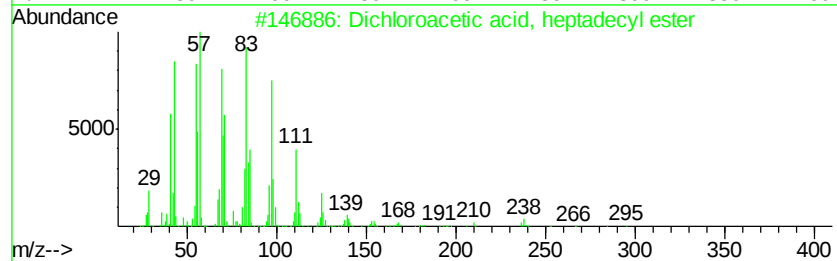
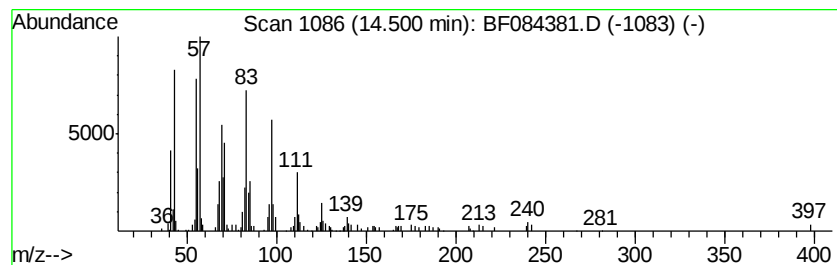
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Dichloroacetic acid, heptadec... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.50	5.58 ng	370030	Chrysene-d12	14.01

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dichloroacetic acid, heptadecyl ...	366	C19H36Cl2O2	1000282-98-2	94
2			Pentafluoropropionic acid, hepta...	402	C20H35F5O2	1000283-04-2	94
3			1-Hexadecanesulfonyl chloride	324	C16H33ClO2S	038775-38-1	91
4			2-Chloropropionic acid, octadec...	360	C21H41ClO2	088104-31-8	91
5			1-Tricosene	322	C23H46	018835-32-0	91



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF011116\
Data File : BF084381.D
Acq On : 11 Jan 2016 15:54
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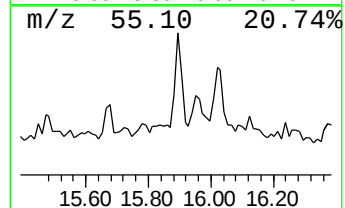
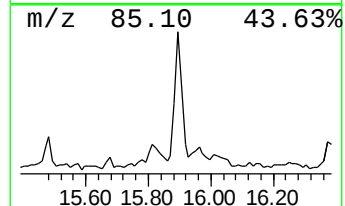
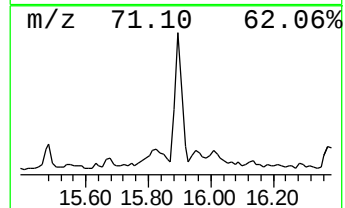
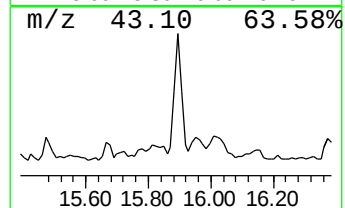
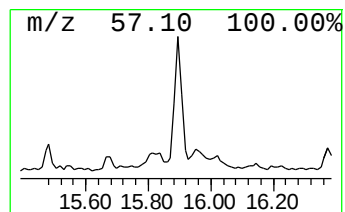
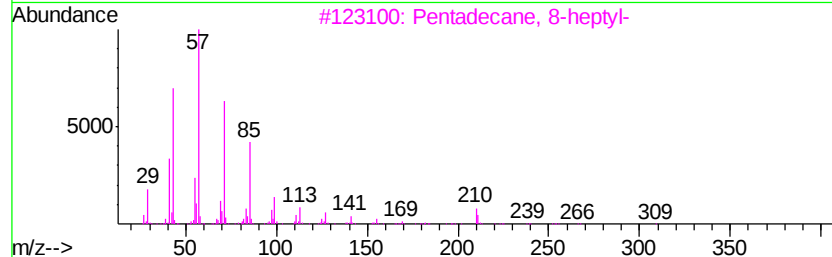
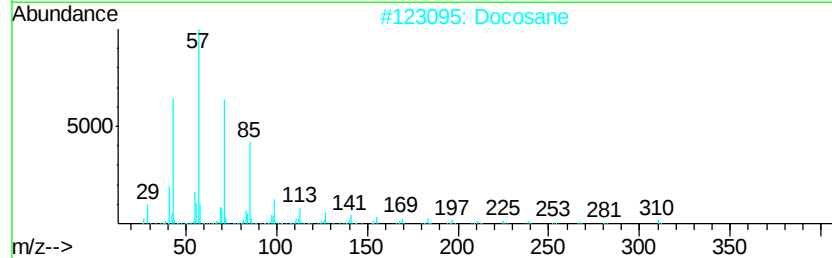
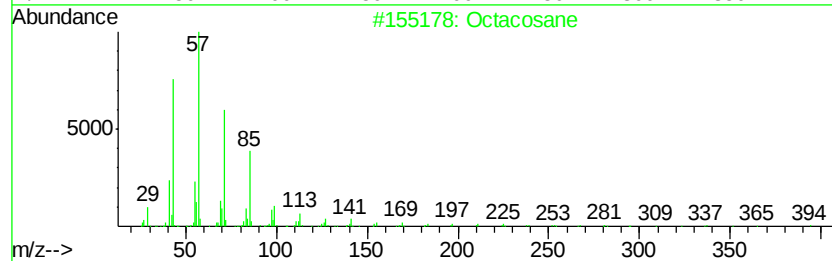
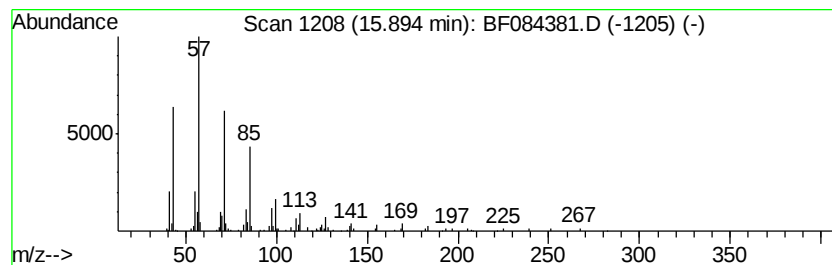
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 Octacosane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.89	3.15 ng	183742	Perylene-d12	15.44

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octacosane	394	C28H58	000630-02-4	91
2		Docosane	310	C22H46	000629-97-0	91
3		Pentadecane, 8-heptyl-	310	C22H46	071005-15-7	91
4		Tetracosane	338	C24H50	000646-31-1	87
5		Tridecane, 1-iodo-	310	C13H27I	035599-77-0	86



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF011116\
Data File : BF084381.D
Acq On : 11 Jan 2016 15:54
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Sample : H1035-02
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ALS Vial : 11 Sample Multiplier: 1

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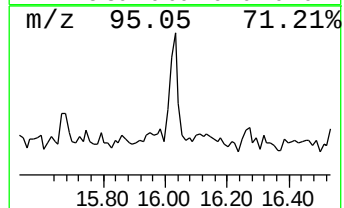
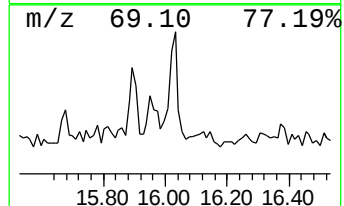
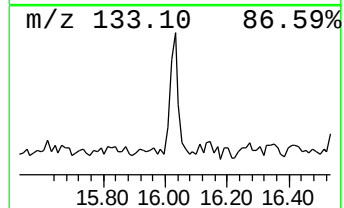
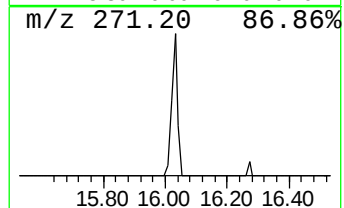
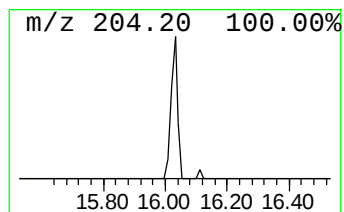
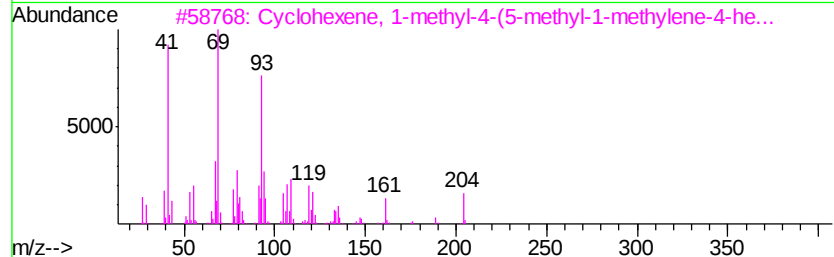
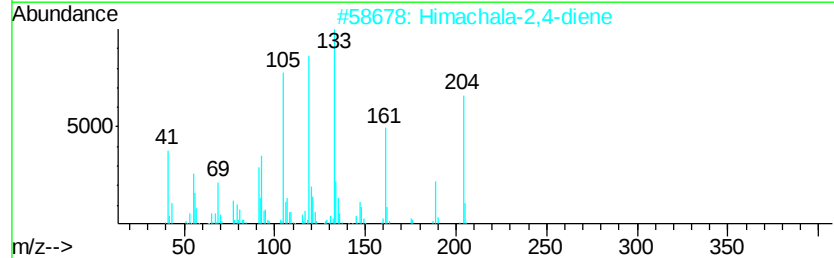
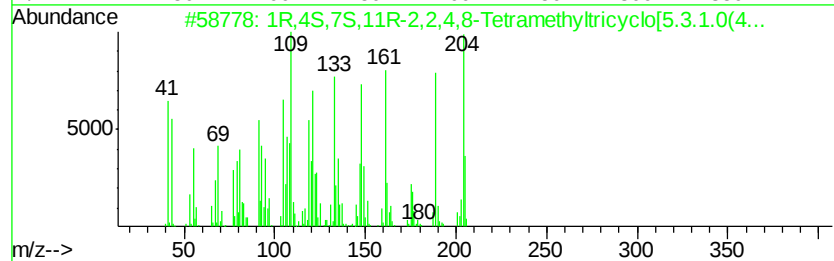
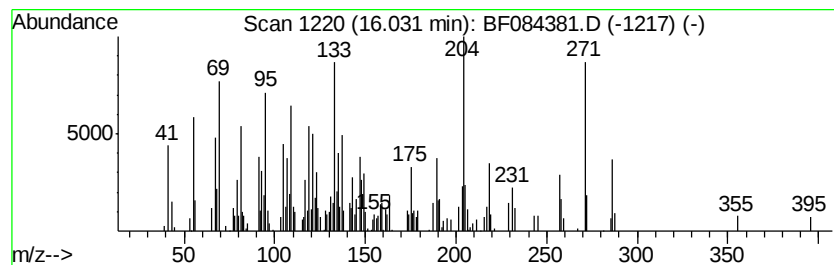
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 unknown16.03 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.03	4.24 ng	247323	Perylene-d12	15.44

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1R,4S,7S,11R-2,2,4,8-Tetramethyl...	204	C15H24	1000140-07-6	35
2		Himachala-2,4-diene	204	C15H24	060909-27-5	30
3		Cyclohexene, 1-methyl-4-(5-methy...	204	C15H24	000495-61-4	27
4		1,2,4,8-Tetramethylbicyclo[6.3.0...	204	C15H24	137235-51-9	22
5		2H-Cyclopentacyclooctene, 4,5,6,...	204	C15H24	1000221-85-8	22



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF011116\
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Misc :
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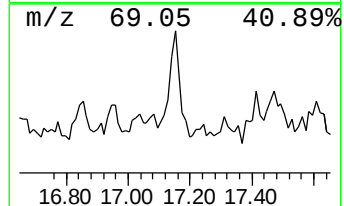
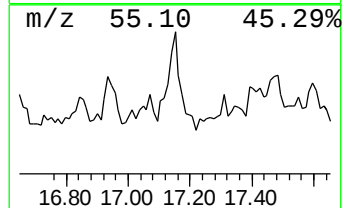
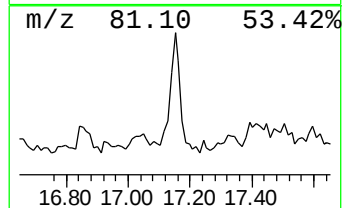
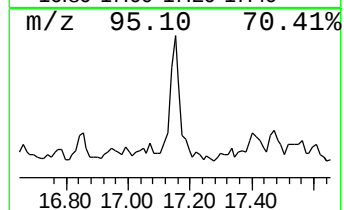
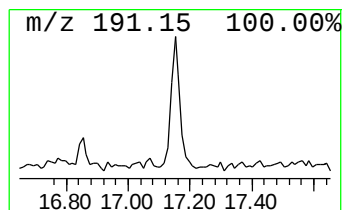
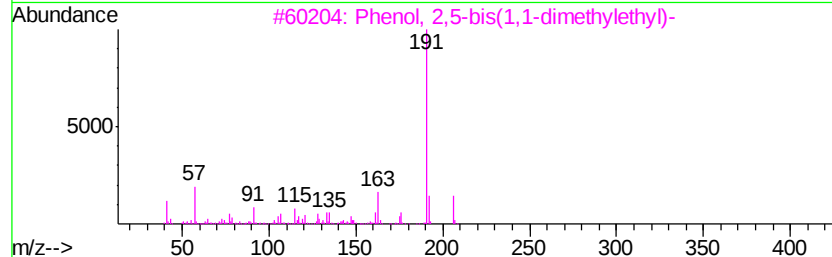
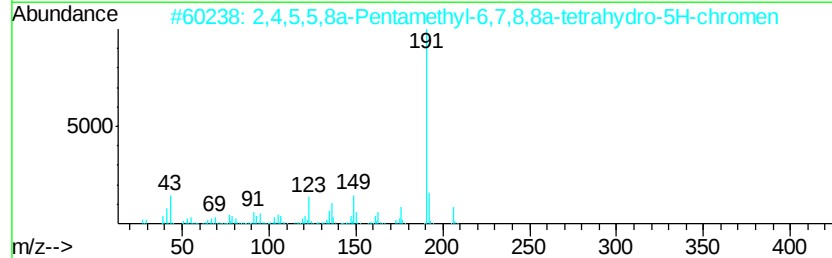
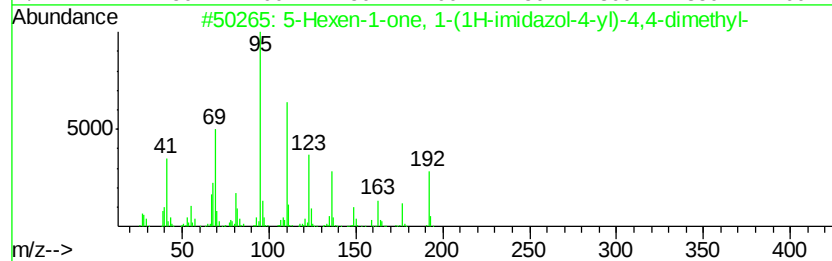
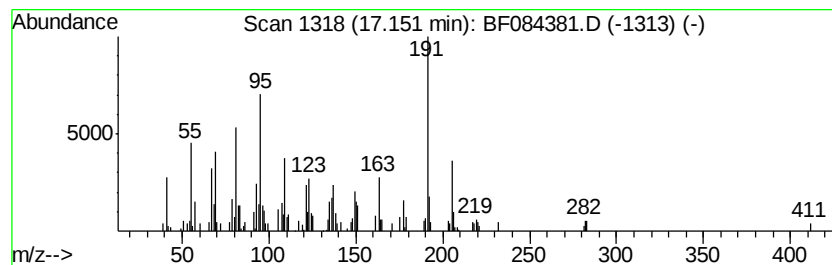
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 18 unknown17.15 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.15	3.70 ng	215700	Perylene-d12	15.44

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Hexen-1-one, 1-(1H-imidazol-4-...	192	C11H16N2O	069393-41-5	47
2		2,4,5,5,8a-Pentamethyl-6,7,8,8a-...	206	C14H22O	1000195-40-9	46
3		Phenol, 2,5-bis(1,1-dimethylethyl)-	206	C14H22O	005875-45-6	38
4		5H-Benzo[def]carbazole	191	C14H9N	000203-65-6	35
5		1-Penten-3-one, 1-(2,6,6-trimeth...	206	C14H22O	000127-43-5	35



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF011116\
Data File : BF084381.D
Acq On : 11 Jan 2016 15:54
Operator : SJ/UM
Sample : H1035-02
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
HAZ-COMP

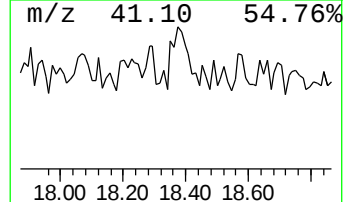
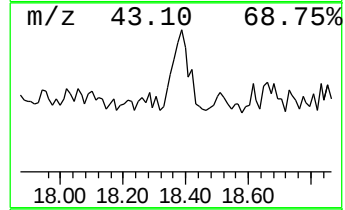
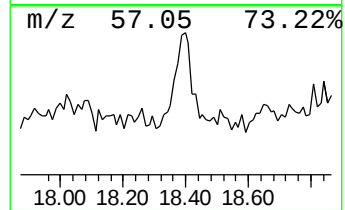
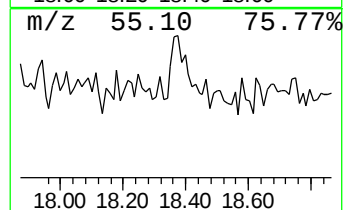
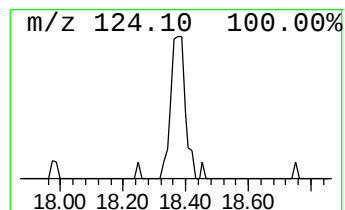
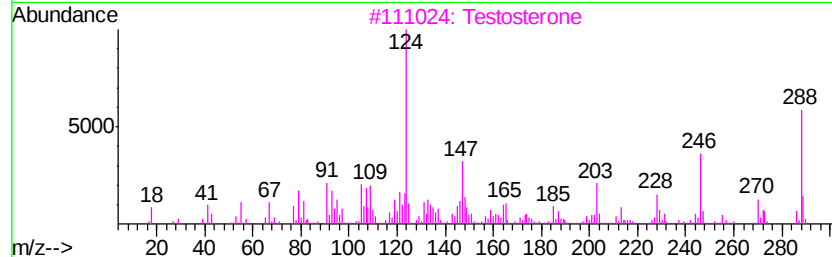
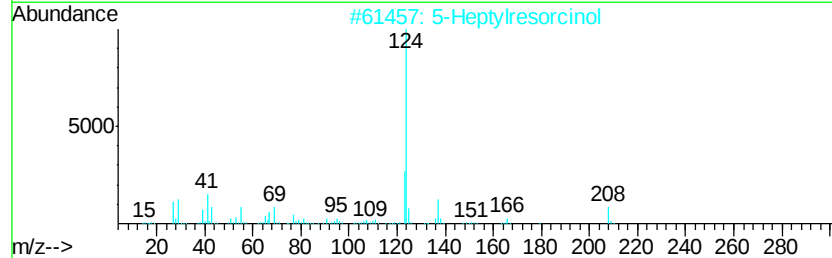
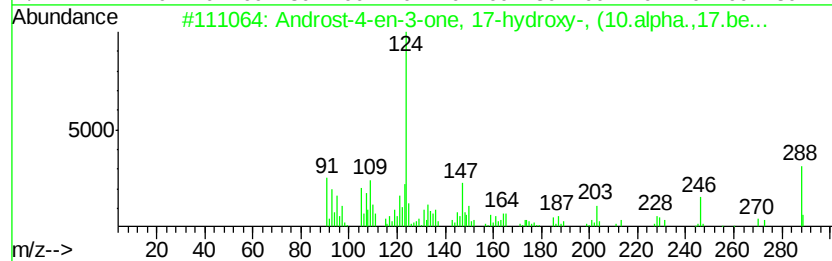
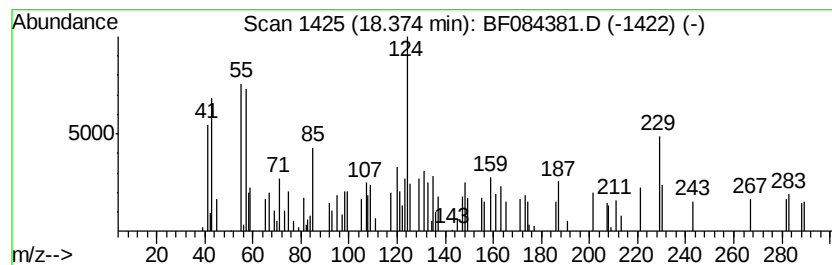
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 20 unknown18.37 Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.37	2.04 ng	118875	Perylene-d12	15.44

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Androst-4-en-3-one, 17-hydroxy-, ...	288	C19H28O2	000604-39-7	42
2		5-Heptylresorcinol	208	C13H20O2	000500-67-4	22
3		Testosterone	288	C19H28O2	000058-22-0	22
4		4H-Pyran-4-one, 2,6-dimethyl-	124	C7H8O2	001004-36-0	14
5		Benzenemethanol, 3-hydroxy-	124	C7H8O2	000620-24-6	14



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF011116\
 Data File : BF084381.D
 Acq On : 11 Jan 2016 15:54
 Operator : SJ/UM
 Sample : H1035-02
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 HAZ-COMP

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF123115.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
2-Pentanone, 4-hy...	5.04	18.4	ng	1484640	1	6.85	1614920	20.0
unknown6.62	6.62	60.0	ng	4846350	1	6.85	1614920	20.0
Naphthalene, 1,2,...	10.02	2.6	ng	284920	3	9.89	2216350	20.0
Eicosane	10.80	2.4	ng	240500	4	11.38	2007000	20.0
1-Perchloroetheny...	12.11	3.1	ng	308948	4	11.38	2007000	20.0
Phenol, 4,4'-(1-m...	12.85	5.4	ng	360293	5	14.01	1326670	20.0
Pentanamide	13.45	20.1	ng	1332500	5	14.01	1326670	20.0
1-Docosene	13.86	10.4	ng	693425	5	14.01	1326670	20.0
Dichloroacetic ac...	14.50	5.6	ng	370030	5	14.01	1326670	20.0
Octacosane	15.89	3.1	ng	183742	6	15.44	1165410	20.0
unknown16.03	16.03	4.2	ng	247323	6	15.44	1165410	20.0
unknown17.15	17.15	3.7	ng	215700	6	15.44	1165410	20.0
unknown18.37	18.37	2.0	ng	118875	6	15.44	1165410	20.0