

Data Path : Z:\HPCHEM1\BNA F\DATA\BF112315\
 Data File : BF083217.D
 Acq On : 23 Nov 2015 21:38
 Operator : UM/IZ
 Sample : G4512-02
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 FILL-DIRT-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-BF112115.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	3.393	116	123	135	rBV	5848002	13530449	100.00%	18.191%
2	4.787	242	245	260	rVB	2037404	3437272	25.40%	4.621%
3	5.244	283	285	288	rBV	4168301	5695652	42.10%	7.658%
4	6.307	375	378	381	rBV	4187986	6359698	47.00%	8.550%
5	6.422	385	388	390	rBV	4271606	5949969	43.97%	8.000%
6	6.639	405	407	409	rBV	1514902	1347036	9.96%	1.811%
7	6.799	418	421	423	rBV	4419755	4537129	33.53%	6.100%
8	7.210	454	457	460	rBV	3336327	3703909	27.37%	4.980%
9	7.930	517	520	524	rBV	1608926	1939506	14.33%	2.608%
10	9.005	611	614	617	rBV	4859405	5610202	41.46%	7.543%
11	9.405	647	649	652	rBV	965548	924517	6.83%	1.243%
12	9.679	670	673	675	rBV	1392747	1807425	13.36%	2.430%
13	10.125	704	712	714	rBV	288864	322696	2.38%	0.434%
14	10.342	728	731	734	rBV	121038	182640	1.35%	0.246%
15	10.468	739	742	744	rBV	2402376	3400508	25.13%	4.572%
16	10.593	751	753	757	rBV	298785	339424	2.51%	0.456%
17	11.039	788	792	794	rBV2	171557	232250	1.72%	0.312%
18	11.154	799	802	805	rVV	1374684	2044332	15.11%	2.749%
19	11.222	805	808	811	rVB2	138857	180718	1.34%	0.243%
20	11.702	848	850	853	rVV	776187	844803	6.24%	1.136%
21	11.759	853	855	861	rVB2	110488	188194	1.39%	0.253%
22	12.354	905	907	909	rBV	737529	673310	4.98%	0.905%
23	12.479	917	918	921	rVV2	214670	317532	2.35%	0.427%
24	12.582	924	927	929	rVV	549321	729611	5.39%	0.981%
25	12.731	938	940	943	rVB	3262169	4300761	31.79%	5.782%
26	12.914	951	956	958	rBV2	81559	161051	1.19%	0.217%
27	13.211	980	982	985	rVB2	116623	147172	1.09%	0.198%
28	13.645	1017	1020	1023	rBV2	384020	477531	3.53%	0.642%
29	13.771	1028	1031	1036	rVV2	1363475	1908992	14.11%	2.567%
30	14.274	1073	1075	1079	rVB2	115009	216639	1.60%	0.291%
31	14.765	1116	1118	1122	rBV	274065	482604	3.57%	0.649%
32	15.028	1139	1141	1144	rVB	171682	215775	1.59%	0.290%
33	15.085	1144	1146	1148	rBV	210266	284439	2.10%	0.382%
34	15.131	1148	1150	1153	rBV	709409	1107998	8.19%	1.490%

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Misc :
ALS Vial : 18 Sample Multiplier: 1

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35	15.554	1185	1187	1189	rBV	122313	201353	1.49%	0.271%
36	16.388	1258	1260	1265	rVB2	97786	202946	1.50%	0.273%
37	16.525	1270	1272	1276	rVB3	114059	223811	1.65%	0.301%
38	16.766	1290	1293	1296	rBV	78457	148896	1.10%	0.200%

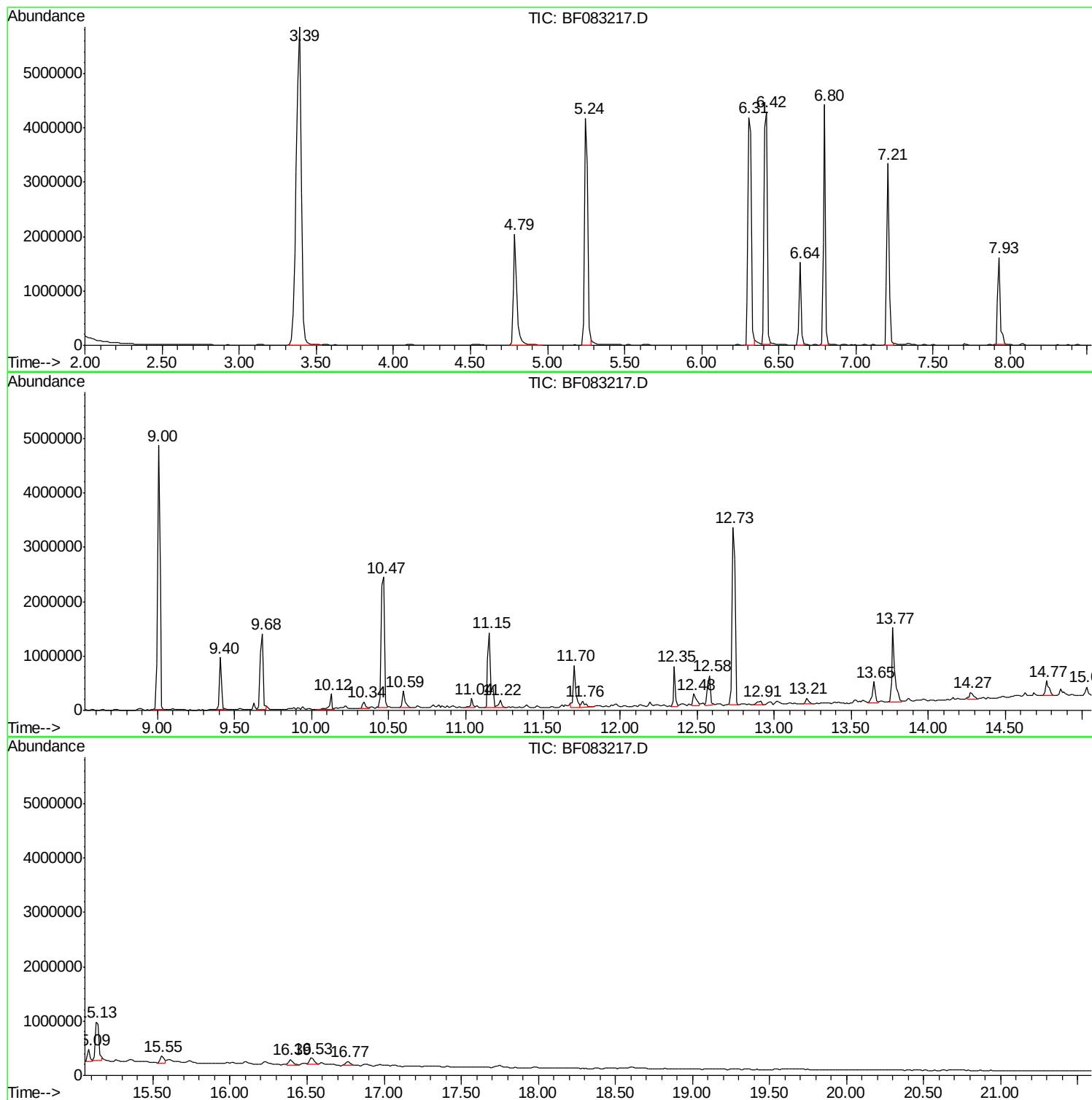
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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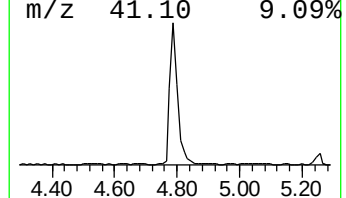
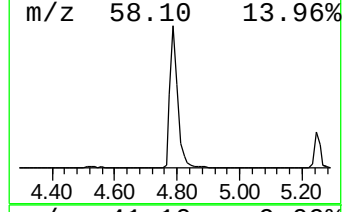
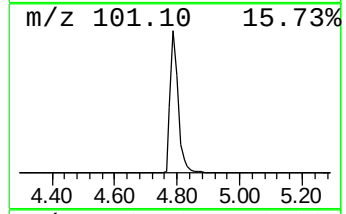
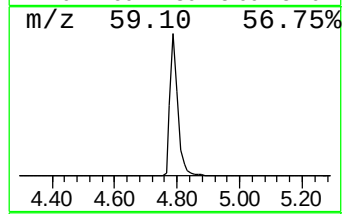
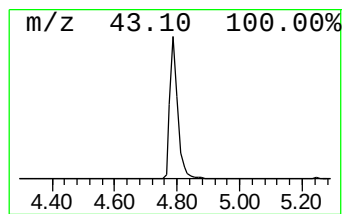
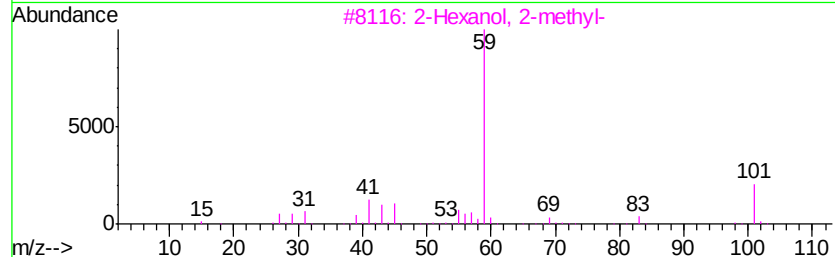
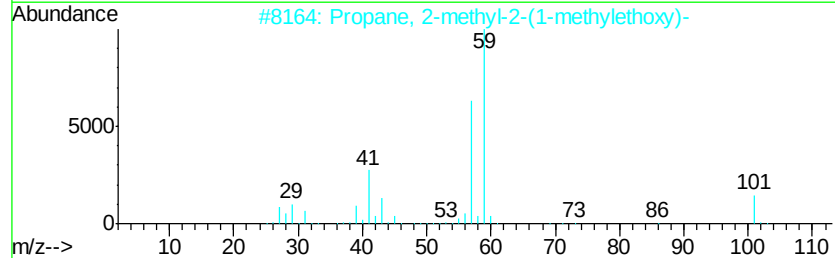
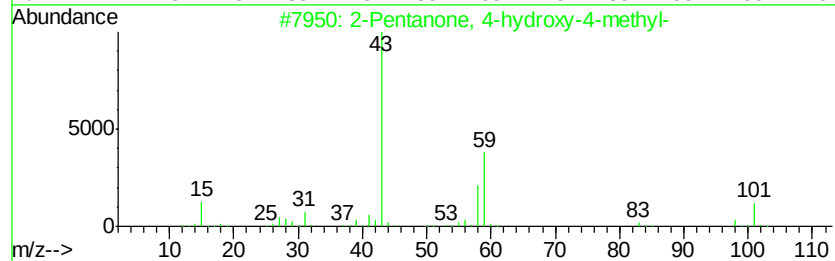
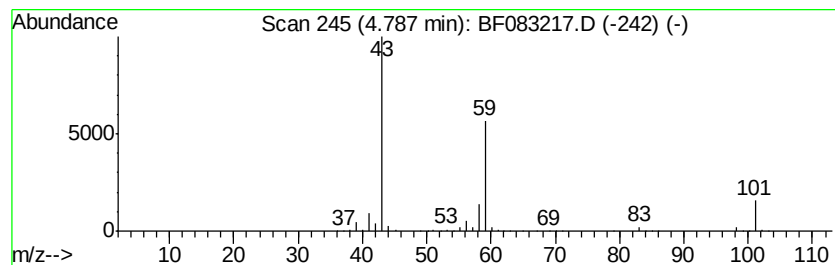
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF112115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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.
4.79	51.03 ng	3437270	1,4-Dichlorobenzene-d4	6.64

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50	
2	Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	42	
3	2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	33	
4	3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33	
5	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	32	



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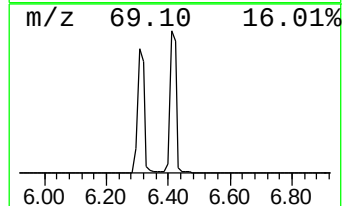
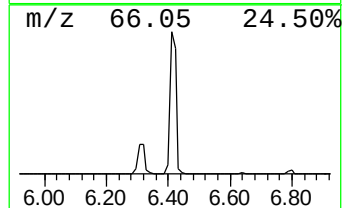
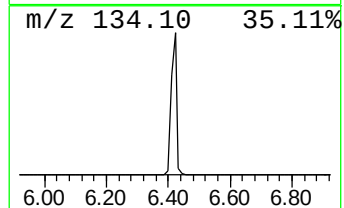
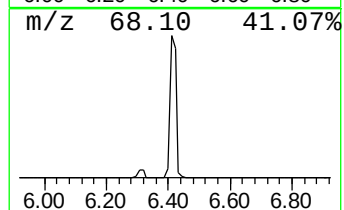
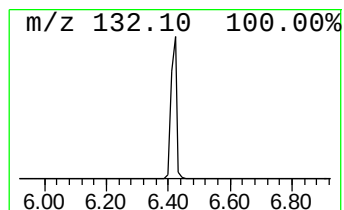
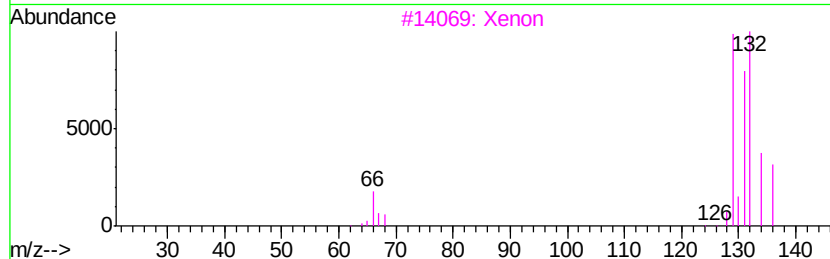
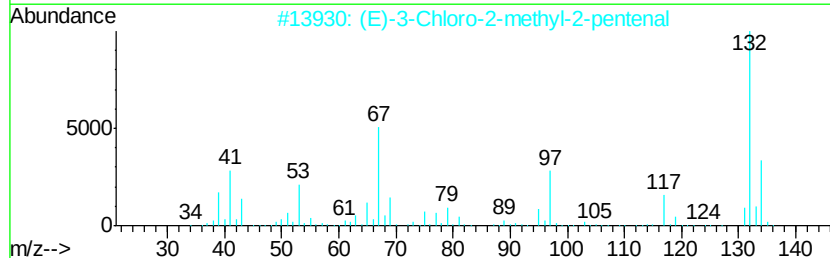
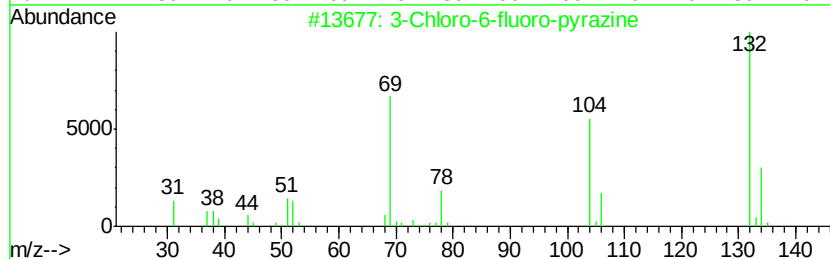
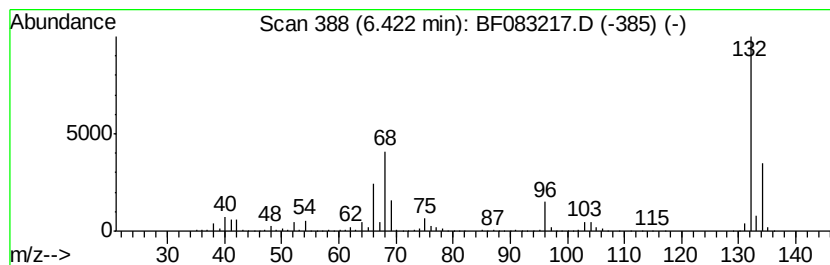
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 3 unknown6.42 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.42	88.34 ng	5949970	1,4-Dichlorobenzene-d4	6.64

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	3	Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2	(E)-3	Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	12
3	Xenon		132	Xe	007440-63-3	9
4	5	Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9
5	1	Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-59-9	9



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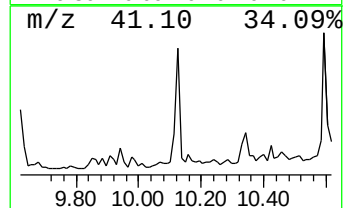
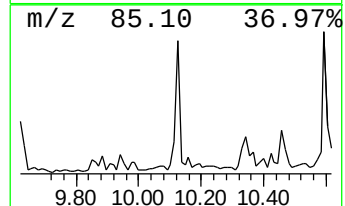
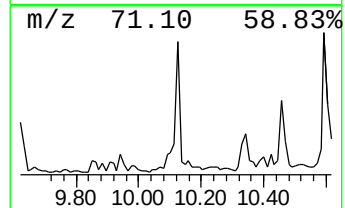
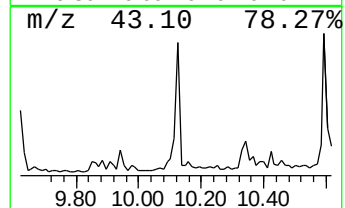
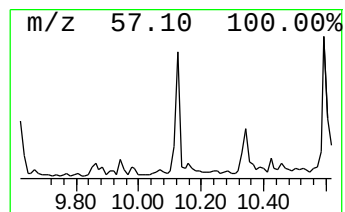
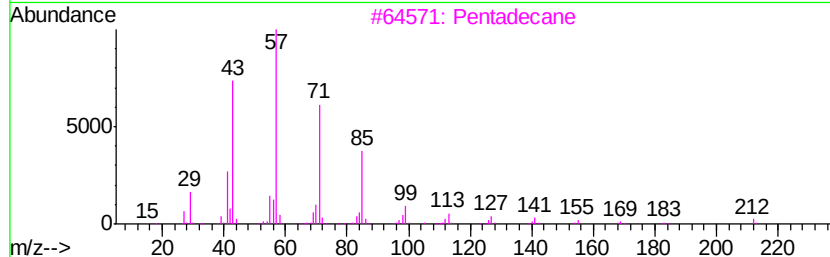
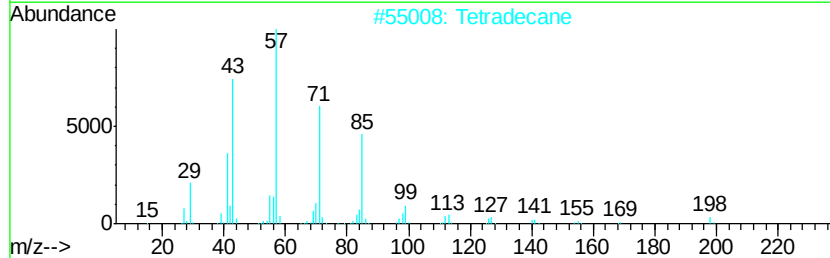
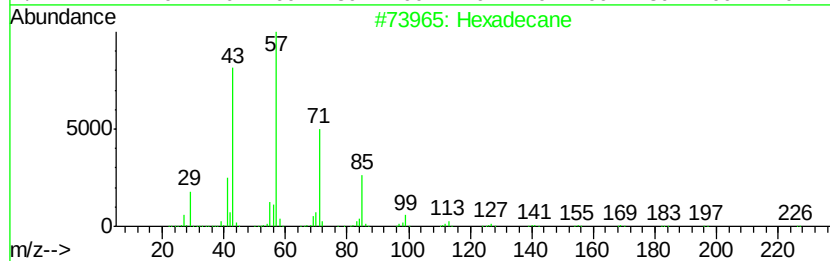
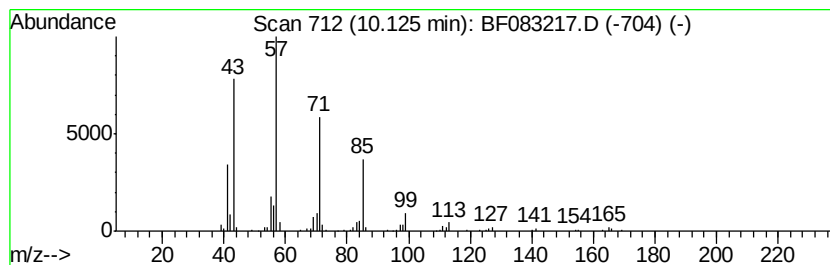
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 5 Hexadecane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.12	3.57 ng	322696	Acenaphthene-d10	9.68

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane		226	C16H34	000544-76-3	91
2	Tetradecane		198	C14H30	000629-59-4	90
3	Pentadecane		212	C15H32	000629-62-9	86
4	Tridecane		184	C13H28	000629-50-5	83
5	Eicosane		282	C20H42	000112-95-8	83



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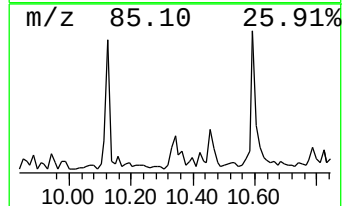
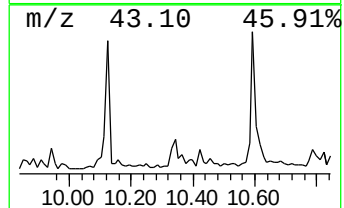
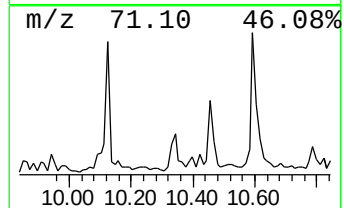
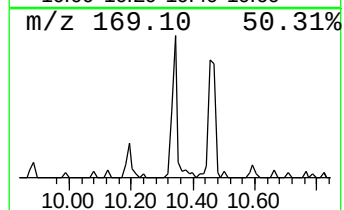
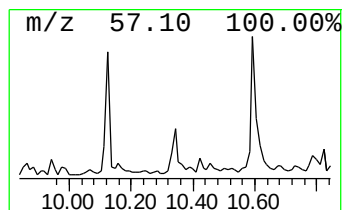
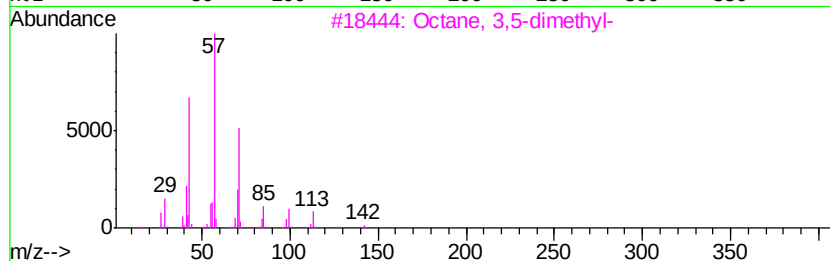
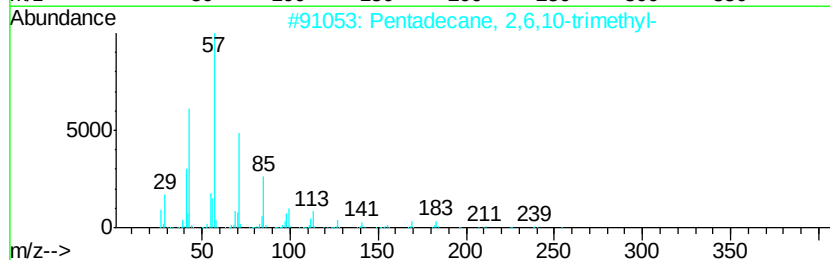
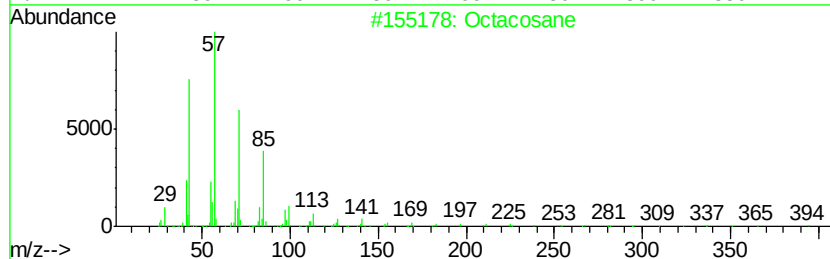
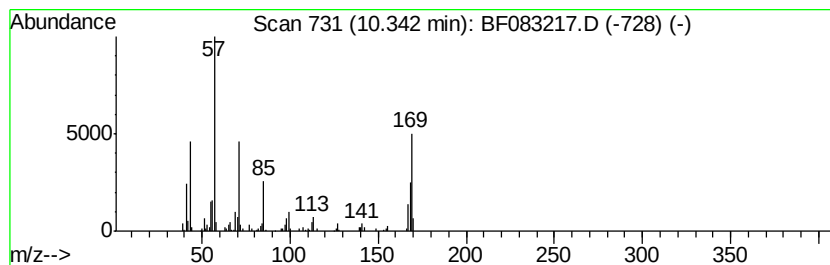
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 unknown10.34 Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.34	2.02 ng	182640	Acenaphthene-d10	9.68

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octacosane	394	C28H58	000630-02-4	46
2		Pentadecane, 2,6,10-trimethyl-	254	C18H38	003892-00-0	46
3		Octane, 3,5-dimethyl-	142	C10H22	015869-93-9	46
4		Heptadecane	240	C17H36	000629-78-7	46
5		Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	43



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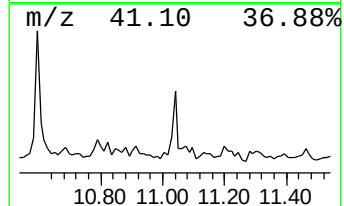
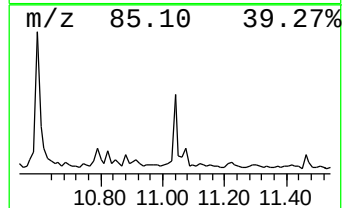
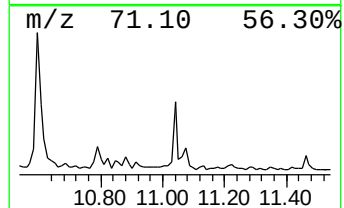
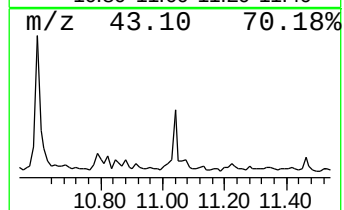
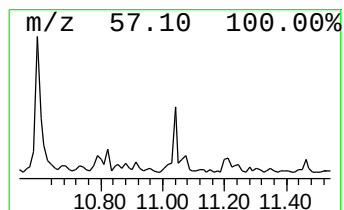
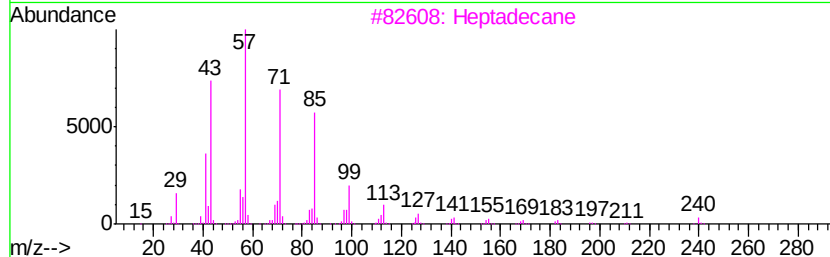
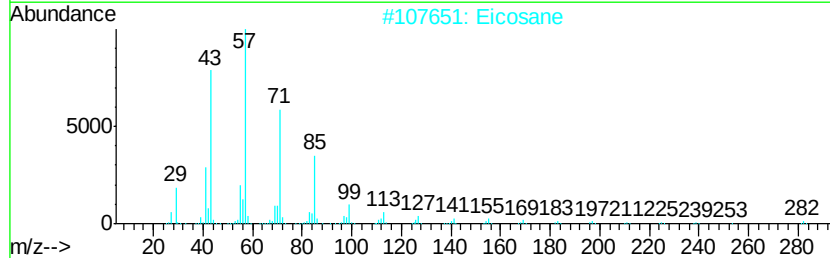
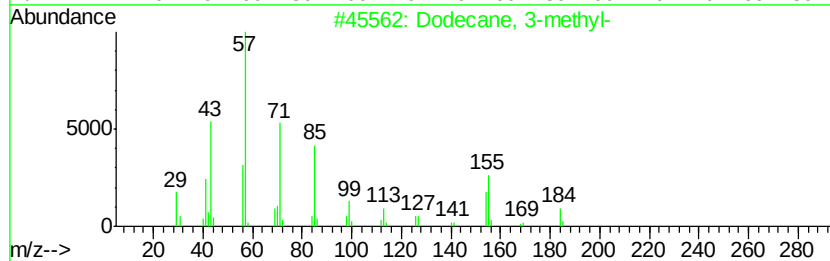
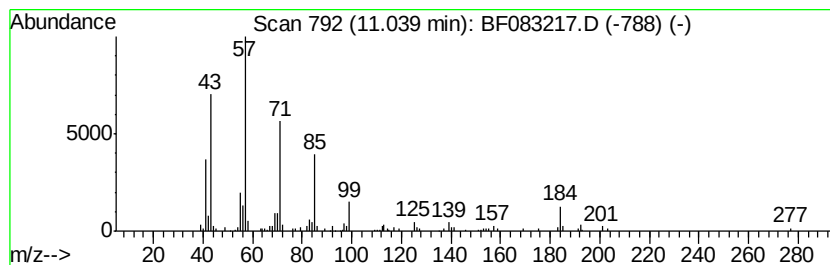
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 8 Dodecane, 3-methyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.04	2.27 ng	232250	Phenanthrene-d10	11.15

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Dodecane, 3-methyl-	184	C13H28	017312-57-1	72
2		Eicosane	282	C20H42	000112-95-8	72
3		Heptadecane	240	C17H36	000629-78-7	64
4		Eicosane, 10-methyl-	296	C21H44	054833-23-7	64
5		Heptacosane	380	C27H56	000593-49-7	64



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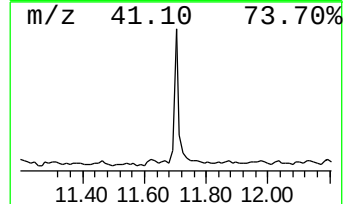
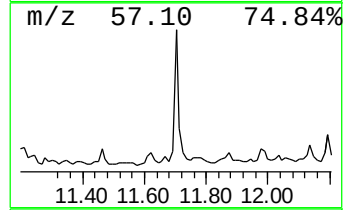
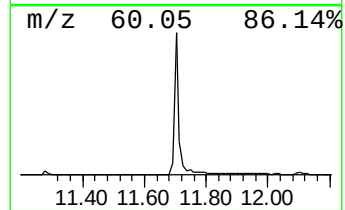
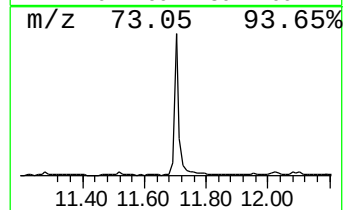
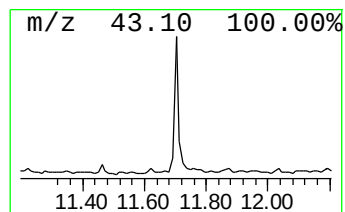
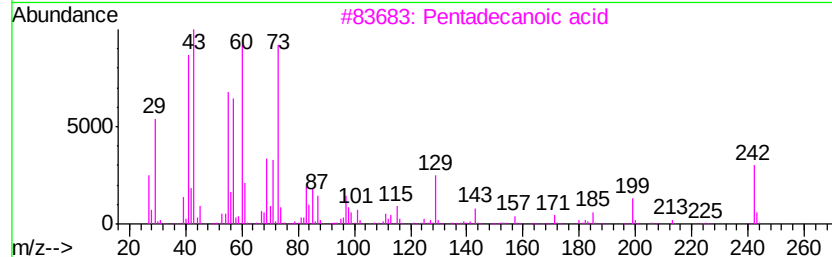
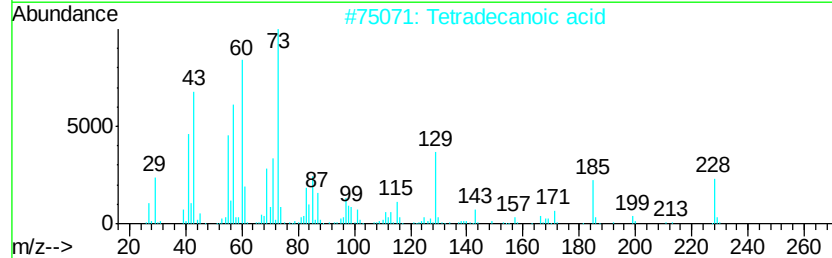
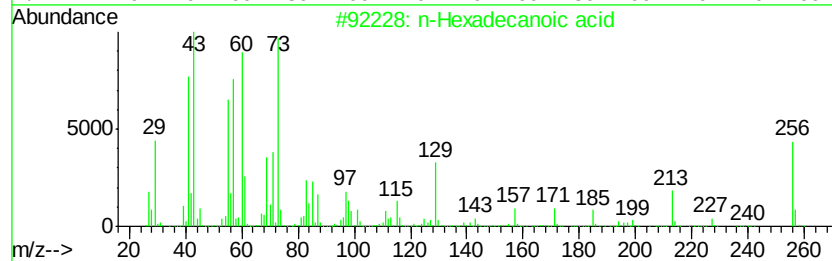
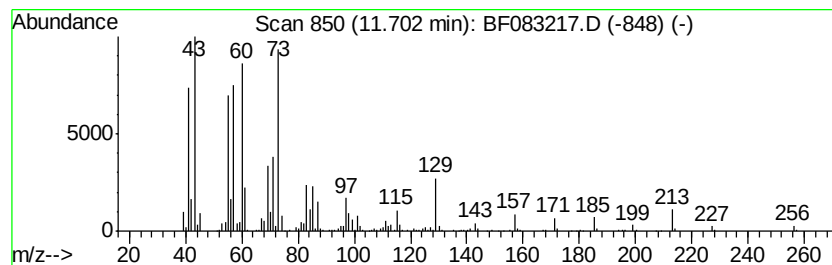
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 n-Hexadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.70	8.26 ng	844803	Phenanthrene-d10	11.15

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2		Tetradecanoic acid	228	C14H28O2	000544-63-8	95
3		Pentadecanoic acid	242	C15H30O2	001002-84-2	87
4		Tridecanoic acid	214	C13H26O2	000638-53-9	80
5		Undecanoic acid	186	C11H22O2	000112-37-8	78



Data Path : Z:\HPCHEM1\BNA F\DATA\BF112315\
Data File : BF083217.D
Acq On : 23 Nov 2015 21:38
Operator : UM/IZ
Sample : G4512-02
Misc :
ALS Vial : 18 Sample Multiplier: 1

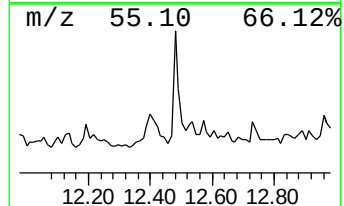
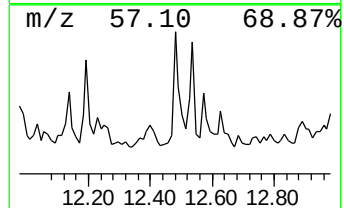
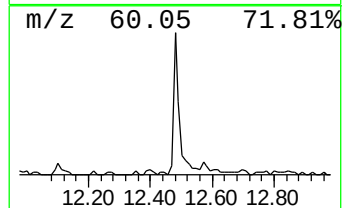
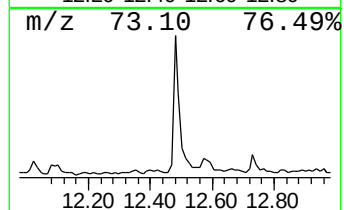
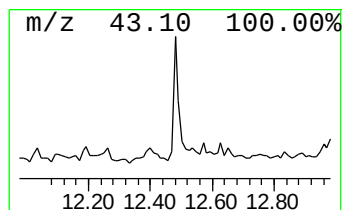
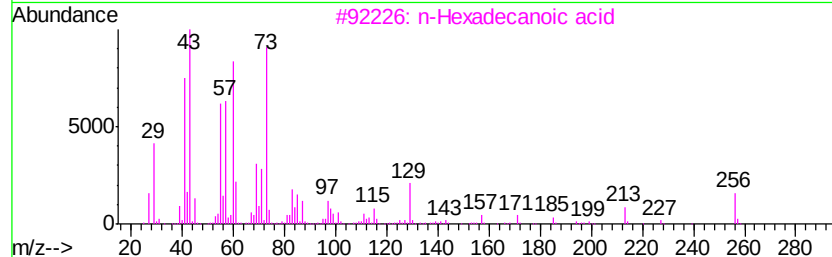
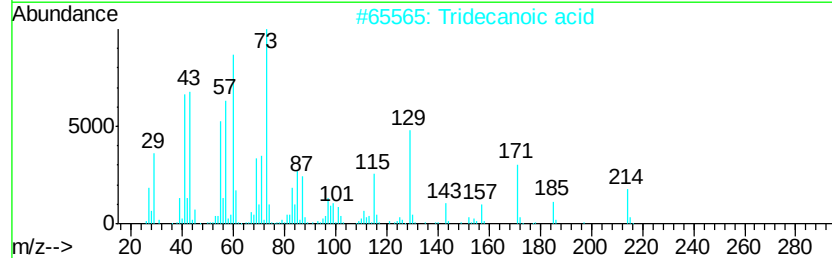
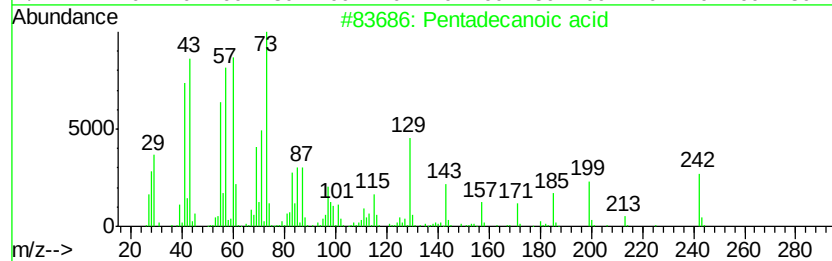
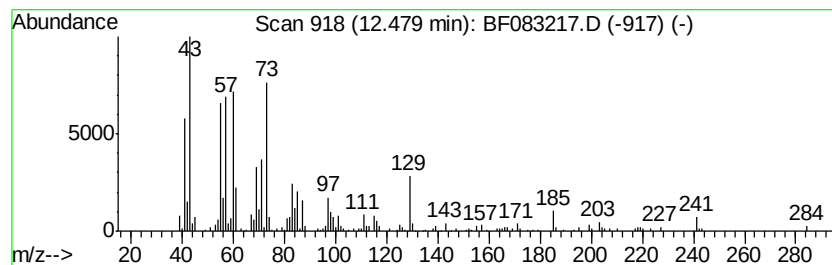
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF112115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Pentadecanoic acid Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD		R.T.
12.48	3.33 ng	317532	Chrysene-d12		13.77
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentadecanoic acid	242	C15H30O2	001002-84-2	94
2	Tridecanoic acid	214	C13H26O2	000638-53-9	78
3	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	53
4	N-Cyclooct-4-enylacetamide	167	C10H17NO	170952-69-9	25
5	Silane, trichlorodocosyl-	442	C22H45Cl3Si	007325-84-0	20



Data Path : Z:\HPCHEM1\BNA F\DATA\BF112315\
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Sample : G4512-02
Misc :
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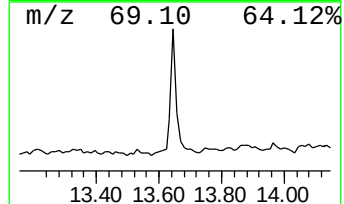
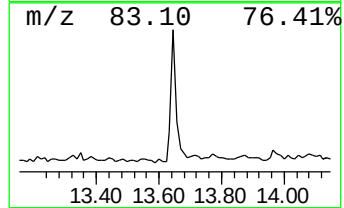
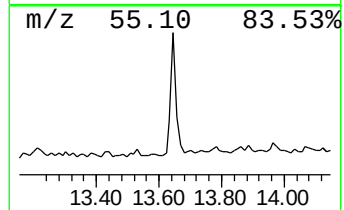
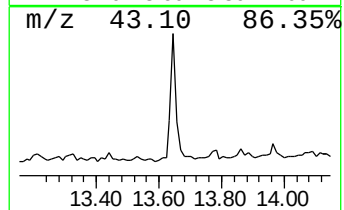
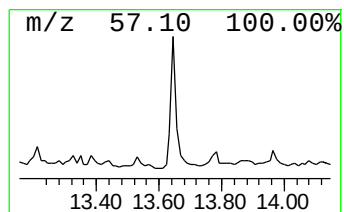
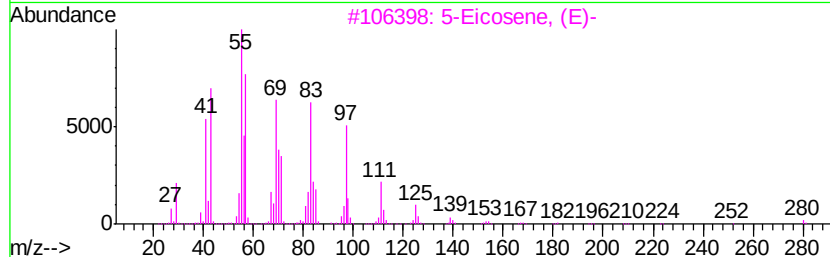
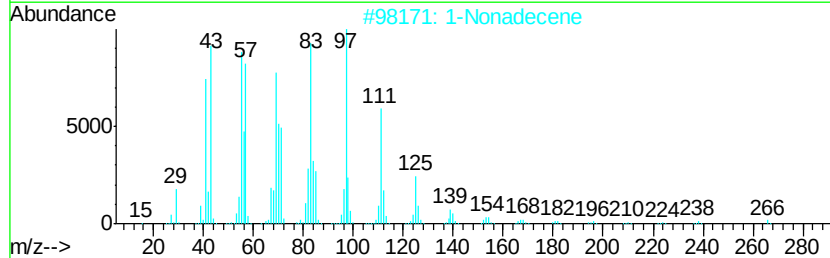
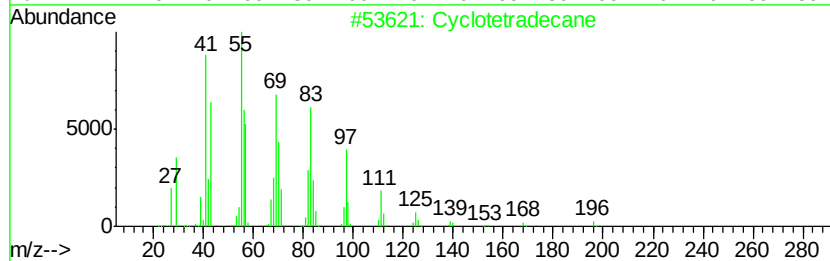
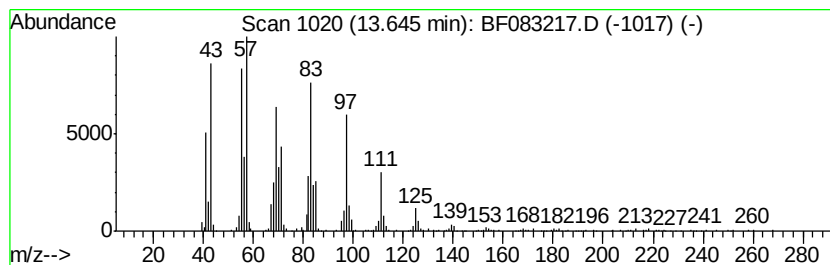
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Cyclotetradecane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.65	5.00 ng	477531	Chrysene-d12	13.77	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclotetradecane	196	C14H28	000295-17-0	93
2	1-Nonadecene	266	C19H38	018435-45-5	91
3	5-Eicosene, (E)-	280	C20H40	074685-30-6	90
4	1-Tetracosanol	354	C24H50O	000506-51-4	87
5	3-Eicosene, (E)-	280	C20H40	074685-33-9	87



Data Path : Z:\HPCHEM1\BNA F\DATA\BF112315\
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Misc :
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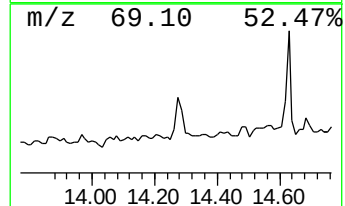
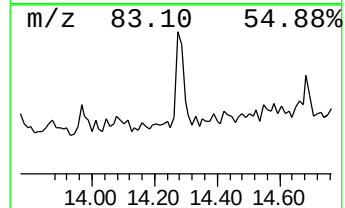
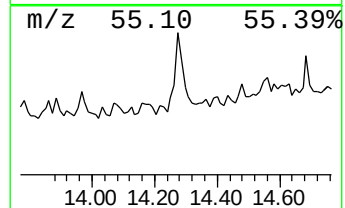
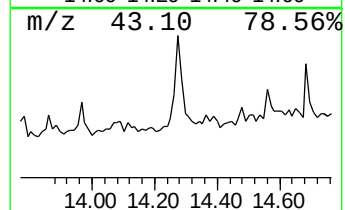
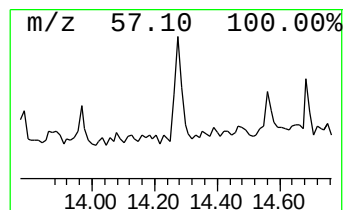
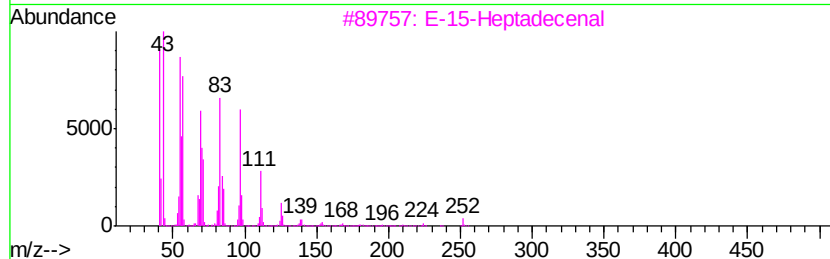
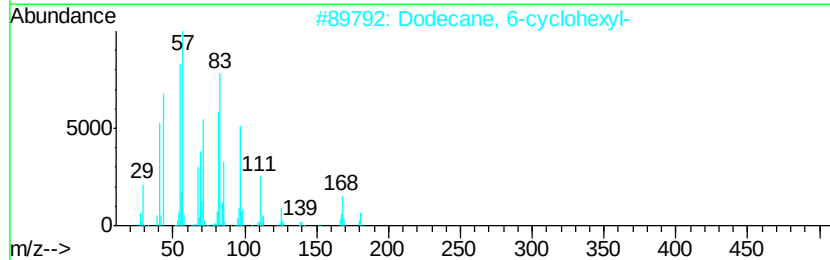
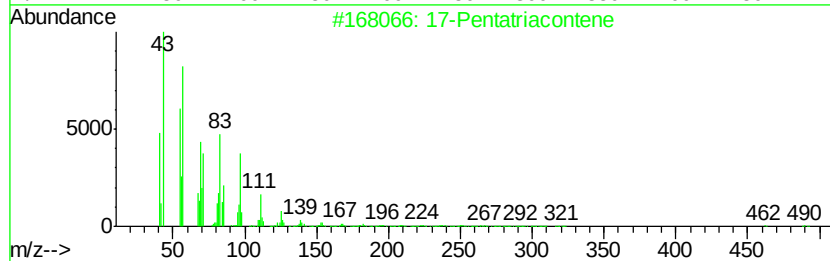
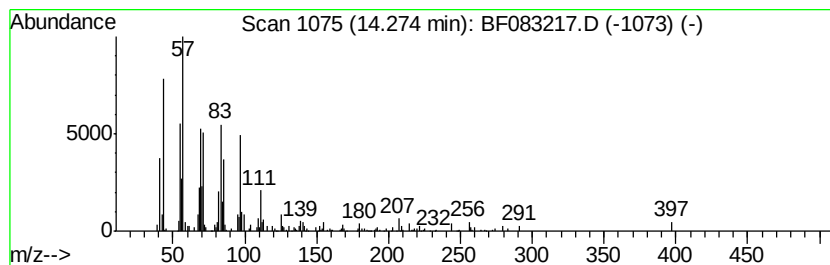
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 17-Pentatriacontene Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.27	2.27 ng	216639	Chrysene-d12	13.77	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	17-Pentatriacontene	491	C35H70	006971-40-0	87
2	Dodecane, 6-cyclohexyl-	252	C18H36	013151-86-5	86
3	E-15-Heptadecenal	252	C17H32O	1000130-97-9	64
4	9-Nonadecene	266	C19H38	031035-07-1	62
5	Cyclotetracosane	336	C24H48	000297-03-0	58



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF112315\
Data File : BF083217.D
Acq On : 23 Nov 2015 21:38
Operator : UM/IZ
Sample : G4512-02
Misc :
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Instrument :
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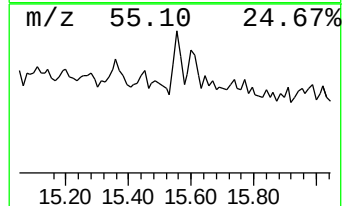
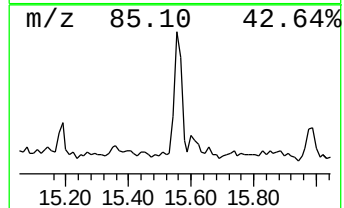
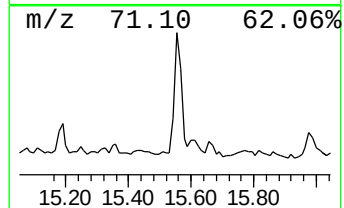
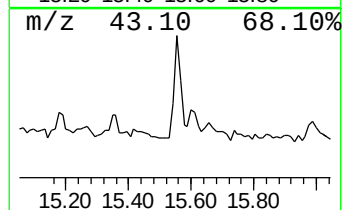
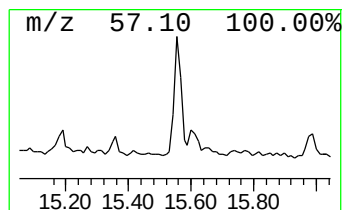
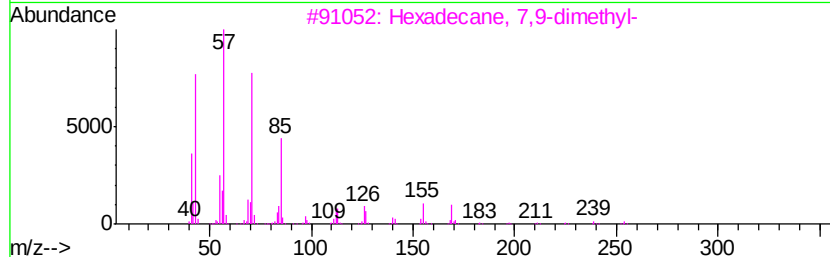
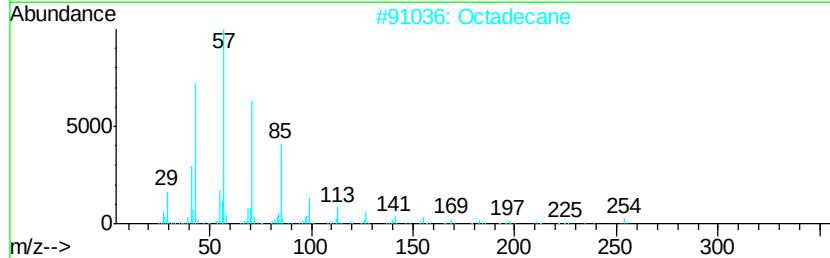
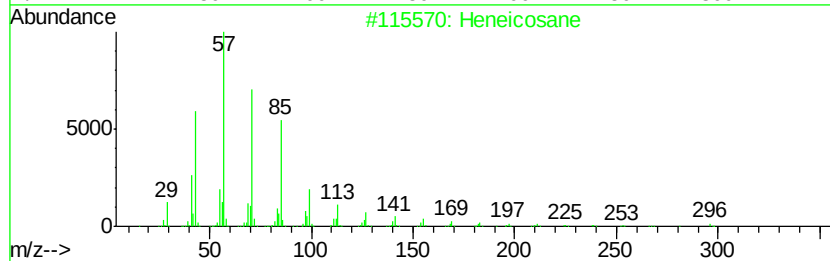
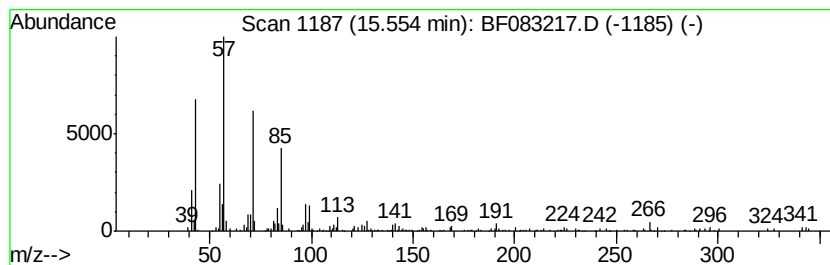
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Heneicosane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.55	3.63 ng	201353	Perylene-d12	15.14

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	91
2		Octadecane	254	C18H38	000593-45-3	90
3		Hexadecane, 7,9-dimethyl-	254	C18H38	021164-95-4	89
4		Tricosane	324	C23H48	000638-67-5	81
5		Octacosane	394	C28H58	000630-02-4	80



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF112315\
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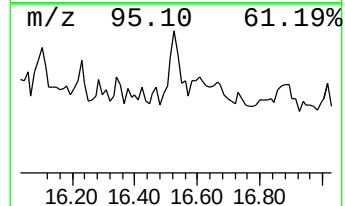
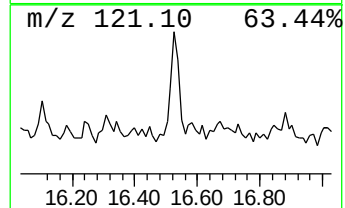
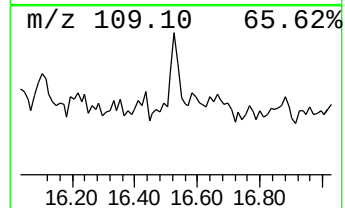
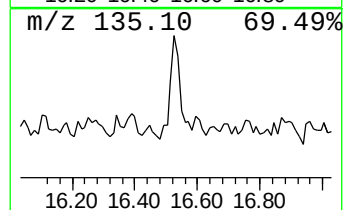
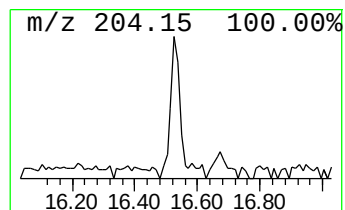
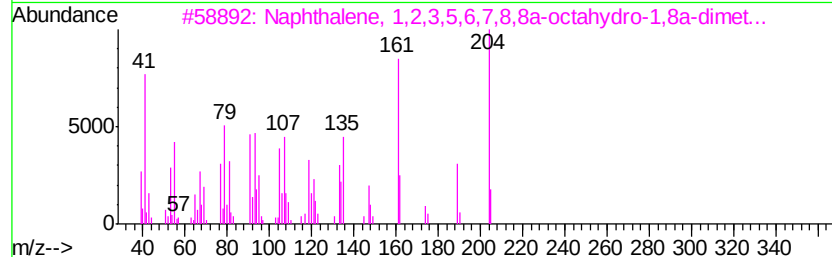
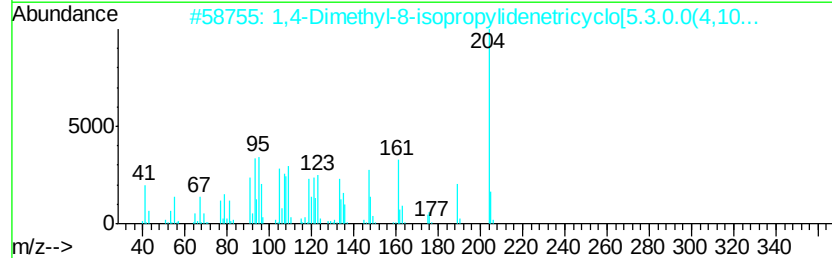
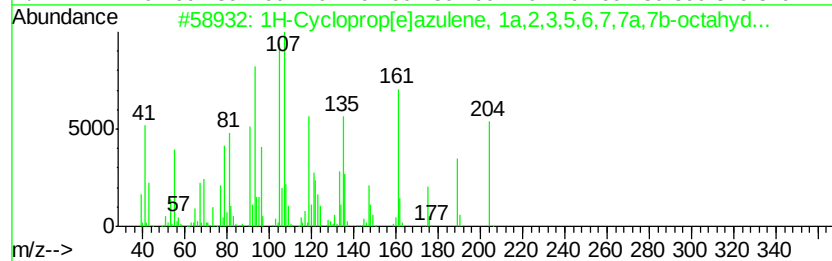
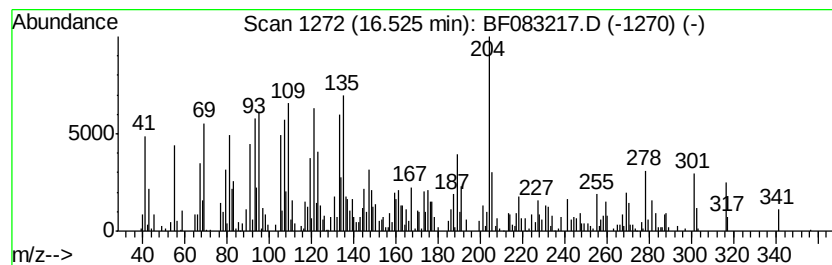
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 1H-Cycloprop[e]azulene, 1a,... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.53	4.04 ng	223811	Perylene-d12	15.14
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		1H-Cycloprop[elazulene, 1a,2,3,5...	204 C15H24	021747-46-6 64
2		1,4-Dimethyl-8-isopropylidenetri...	204 C15H24	1000140-07-7 64
3		Naphthalene, 1,2,3,5,6,7,8,8a-oc...	204 C15H24	004630-07-3 47
4		Aciphyllene	204 C15H24	087745-31-1 41
5		1R,4S,7S,11R-2,2,4,8-Tetramethyl...	204 C15H24	1000140-07-6 35



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF112315\
Data File : BF083217.D
Acq On : 23 Nov 2015 21:38
Operator : UM/IZ
Sample : G4512-02
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF112115.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
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unknown6.42	6.42	88.3	ng	5949970	1	6.64	1347040	20.0
Hexadecane	10.12	3.6	ng	322696	3	9.68	1807430	20.0
unknown10.34	10.34	2.0	ng	182640	3	9.68	1807430	20.0
Dodecane, 3-methyl-	11.04	2.3	ng	232250	4	11.15	2044330	20.0
n-Hexadecanoic acid	11.70	8.3	ng	844803	4	11.15	2044330	20.0
Pentadecanoic acid	12.48	3.3	ng	317532	5	13.77	1908990	20.0
Cyclotetradecane	13.65	5.0	ng	477531	5	13.77	1908990	20.0
17-Pentatriacontene	14.27	2.3	ng	216639	5	13.77	1908990	20.0
Heneicosane	15.55	3.6	ng	201353	6	15.14	1108000	20.0
1H-Cycloprop[e]az...	16.53	4.0	ng	223811	6	15.14	1108000	20.0