

# Package ‘BayesX’

February 19, 2015

**Type** Package

**Title** R Utilities Accompanying the Software Package BayesX

**Version** 0.2-9

**Date** 2014-08-18

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**Description** This package provides functionality for exploring and visualising estimation results obtained with the software package BayesX for structured additive regression. It also provides functions that allow to read, write and manipulate map objects that are required in spatial analyses performed with BayesX, a free software for estimating structured additive regression models (<http://www.bayesx.org>).

**Depends** shapefiles

**Imports** sp, maptools, colorspace (>= 0.97), coda, splines

**Suggests** spdep, akima, gplotlib, rgeos

**LazyLoad** yes

**License** GPL-2

**NeedsCompilation** no

**Repository** CRAN

**Date/Publication** 2014-08-18 16:46:47

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BayesX-package	<i>R Utilities Accompanying the Software Package BayesX</i>
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## Description

This package provides functionality for exploring and visualising estimation results obtained with the software package BayesX for structured additive regression. It also provides functions that allow to read, write and manipulate map objects that are required in spatial analyses performed with BayesX.

## Details

Package: BayesX  
 Type: Package  
 Version: 0.2-6  
 Date: 2013-03-11  
 License: GPL-2

## Author(s)

Thomas Kneib, Felix Heinzl, Andreas Brezger, Daniel Sabanes Bove  
 Maintainer: Thomas Kneib <tkneib@uni-goettingen.de>

## References

<http://www.bayesx.org>

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add.neighbor	<i>Add Neighborhood Relations</i>
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## Description

Adds a neighborhood relationship between two given regions to a map object in graph format.

## Usage

```
add.neighbor(map, region1, region2)
```

## Arguments

map	Map object in graph format that should be modified.
region1, region2	Names of the regions that should be connected as neighbors.

## Value

Returns an adjacency matrix that represents the neighborhood structure of map plus the new neighborhood relation in graph format.

## Author(s)

Felix Heinzl, Thomas Kneib

## See Also

[get.neighbor](#), [delete.neighbor](#), [read.gra](#), [write.grabnd2gra](#)

## Examples

```
germany <- read.gra(system.file("examples/germany.gra", package="BayesX"))
get.neighbor(germany, c("1001", "7339"))
germany <- add.neighbor(germany, "7339", "1001")
get.neighbor(germany, c("1001", "7339"))
```

---

`bnd2gra`*Convert Boundary Format to Graph Format*

---

**Description**

Converts a map in boundary format to a map in graph format.

**Usage**

```
bnd2gra(map)
```

**Arguments**

`map` Map in boundary format that should be converted.

**Value**

Returns an adjacency matrix that represents the neighborhood structure of the map object in graph format.

**Author(s)**

Felix Heinzl, Thomas Kneib

**References**

BayesX Reference Manual. Available from <http://www.bayesx.org>

**See Also**

[read.bnd](#), [read.gra](#), [write.bnd](#), [write.gra](#)

**Examples**

```
tanzania.bnd <- read.bnd(system.file("examples/tanzania.bnd", package="BayesX"))
tanzania.gra <- bnd2gra(tanzania.bnd)
```

---

`createxymap`*Create map objects for some points and a given distance*

---

### Description

Creates a map object from a list of coordinates by treating observations within a certain distance as neighbors. The resulting neighborhood structure is stored in a map object in graph format while a map in boundary format is created to enable visualisation.

### Usage

```
createxymap(x, y, districts=NULL, p=2, max.dist)
```

### Arguments

<code>x</code>	Vector of x-coordinates of underlying points
<code>y</code>	Vector of y-coordinates of underlying points
<code>districts</code>	Either NULL or a vector of names for labeling points. If <code>districts=NULL</code> , points are labelled by index.
<code>p</code>	Any p-norm with $p \geq 1$ can be chosen as the distance measure with the Euclidian distance ( $p=2$ ) being the default. For $p=Inf$ , the maximum of coordinates is used. Except for $p=Inf$ , calculations can fail for huge $p$ .
<code>max.dist</code>	Value which determines the neighborhood. Points whose distance is smaller or equal than <code>max.dist</code> are considered as neighbors.

### Value

List of two elements: map object in graph format and map object in boundary format.

### Author(s)

Felix Heinzl, Thomas Kneib

### See Also

[read.gra](#), [read.bnd](#), [drawmap](#)

### Examples

```
x <- c(3,3,5,9.5,11,11)
y <- c(2,4,4,6,4.5,5)
xymap <- createxymap(x, y, districts=c("A","B","C","D","E","F"), max.dist=2)
xymap$gra
drawmap(map=xymap$bnd)
```

---

delete.neighbor	<i>Delete Neighborhood Relations</i>
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### Description

Adds the neighborhood relationship between two given regions from a map object in graph format.

### Usage

```
delete.neighbor(map, region1, region2)
```

### Arguments

map	Map object in graph format that should be modified.
region1, region2	Names of the regions that should no longer be regarded as neighbors.

### Value

Returns an adjacency matrix that represents the neighborhood structure of map minus the deleted neighborhood relation in graph format.

### Author(s)

Felix Heinzl, Thomas Kneib

### See Also

[get.neighbor](#), [add.neighbor](#), [read.gra](#), [write.grabnd2gra](#)

### Examples

```
germany <- read.gra(system.file("examples/germany.gra", package="BayesX"))
get.neighbor(germany, c("7339"))
germany <- delete.neighbor(germany, "7339", "7141")
get.neighbor(germany, c("7339"))
```

**Description**

Visualises variables that are spatially aligned according to a given map object. Each of the regions in a map will be coloured according to the value of the variable.

**Usage**

```
drawmap(data, map, regionvar=2, plotvar=3, limits, cols="hcl", nrcolors=100,
        swapcolors=FALSE, pcat=FALSE,
        hcl.par=list(h=c(120, 0), c=60, l=c(45,60), power=1.2),
        hsv.par=list(s=1, v=1), legend=TRUE, drawnames=FALSE, cex.names=0.7,
        cex.legend=0.7, mar.min=2, density=15, ...)
```

**Arguments**

data	Either the name of a file or a data frame containing the variables to be visualised. If missing, the map will be visualised without superposition of any further information
map	Map object containing the required boundary information (as obtained by a call to <a href="#">read.bnd</a> )
regionvar	Defines the variable specifying the geographical regions. Either the name of a variable in data or the index of the corresponding column.
plotvar	Defines the variable that should be visualised. Either the name of a variable in data or the index of the corresponding column.
limits	Restricts (or extends) the coloring scheme to a range of values.
cols	Color scheme to be employed. Could be either a vector of colors or one out of the following pre-defined schemes: hcl, hsv, grey
nrcolors	Number of colors (only meaningful when using one of the pre-defined colour schemes).
swapcolors	Reverse the order of colors (works also with user-specified colours but will be most useful with the pre-defined schemes).
pcat	Option for the visualisation of posterior probabilities. In this case, a three-colour scheme representing significantly positive, insignificant and significantly negative values.
hcl.par	Parameters for the hcl colour scheme (see the documentation of <code>diverge_hcl</code> in package <code>vcd</code> for details).
hsv.par	Parameters for the hsv colour scheme (see the documentation of <code>hsv</code> for details).
legend	Should a legend be added to the figure?
drawnames	Adds the name of each region as a text label to the plot. In most cases the result will be confusing but may be useful when checking the validity of a map.

<code>cex.names</code>	Magnification to be used for the names (if <code>drawnames=TRUE</code> ).
<code>cex.legend</code>	Magnification to be used for the legend.
<code>mar.min</code>	Controls the definition of boundaries. Could be either <code>NULL</code> for individual settings of <code>mar</code> or a value which defines <code>mar</code> as follows: The boundaries will be calculated according to the height to width ratio of the map with minimal boundary <code>mar.min</code> .
<code>density</code>	Regions without data will be visualised with diagonal stripes. <code>density</code> defines how dense the stripes should be.
<code>...</code>	Further arguments to be passed to the plot calls that visualise the region boundaries (probably not useful at all).

**Author(s)**

Felix Heinzl, Thomas Kneib, Andreas Brezger

**See Also**

[read.bnd](#)

**Examples**

```

germany <- read.bnd(system.file("examples/germany.bnd", package="BayesX"))
drawmap(map=germany)
drawmap(map=germany, drawnames=TRUE)

res <- read.table(system.file("examples/spatial_f_regions_spatial.res",
                             package="BayesX"), header=TRUE)

drawmap(res, map=germany)
drawmap(res, map=germany, limits=c(-2,4))
drawmap(res, map=germany, regionvar="regions", plotvar="pmed")
drawmap(res, map=germany, legend=FALSE)
drawmap(res, map=germany, legend=FALSE, main="spatial effect")

drawmap(res, map=germany, cols="hsv")
drawmap(res, map=germany, swapcolors=TRUE, cols="hsv")
drawmap(res, map=germany, cols="grey")
drawmap(res, map=germany,
        cols=c('darkgreen', 'green', 'yellow', 'orange', 'red', 'darkred'))

drawmap(res, map=germany, pcat=TRUE, cols="hcl")
drawmap(res, map=germany, pcat=TRUE, cols="hsv")
drawmap(res, map=germany, pcat=TRUE, cols="grey")

drawmap(res, map=germany, nrcolors=10, cols="hcl")
drawmap(res, map=germany, nrcolors=10, cols="hsv")
drawmap(res, map=germany, nrcolors=10, cols="grey")

drawmap(res, map=germany, cols="hcl",
        hcl.par=list(h=c(0,120), c=60, l=c(45,90), power=1.2))

```



```
drawmap(res, map=germany, cols="hcl",
        hcl.par=list(h=c(300,120), c=60, l=c(45,90), power=1.2))
drawmap(res, map=germany, cols="hcl",
        hcl.par=list(h=c(40,260), c=60, l=c(45,90), power=1.2))
drawmap(res, map=germany, cols="hsv", hsv.par=list(s=0.7, v=0.7))
```

---

extractSamples	<i>Extract MCMC samples from a BayesX results directory</i>
----------------	---

---

### Description

This is a convenience function to extract samples from a BayesX results directory, which processes the log file to e.g. convert the spline coefficients samples to function values samples.

### Usage

```
extractSamples(directoryWithBasename,
              logfile = file.path(dirname(directoryWithBasename), "log.txt"))
```

### Arguments

directoryWithBasename	The BayesX results directory with basename for the files (e.g. "results/test", if this was specified as outfile in BayesX for the bayesreg object)
logfile	The log file of the MCMC run, defaults to log.txt in the results directory.

### Value

Returns a list with the extracted samples of effects and deviances as well as the prediction data.frame:

<function name>

for P-Splines, Random Walks and spatial effects: a list with mcmc objects 'functionSamples' and 'varianceSamples' containing the respective effects/function and variance parameter samples.

FixedEffects an mcmc object of all fixed simple parametric effects

RandomEffects if there is at least one random effect in the model, this is a list, with elements in the first hierarchy being the group ID names, and elements in the second hierarchy being the names of the covariates. The leafs are the mcmc objects 'functionSamples' and 'varianceSamples', as for the other non-fixed terms

Deviance an mcmc object with the (unstandardized and saturated) deviance

means if the option predictmu was used, this mcmc object contains the mean samples

scale an mcmc object with the possible scale parameter samples

lassoCoefficients

an mcmc object with the possible lasso regression parameter samples

ridgeCoefficients

an mcmc object with the possible ridge regression parameter samples

PredictMeans data.frame corresponding to the possible predictmean file in the BayesX directory

Additionally, entries for possibly remaining lasso or ridge variance parameters etc. are included in the return list.

### Warning

You should be sure that only one MCMC run is saved in the given results directory in order to get sensible results out of this function.

### Author(s)

Daniel Sabanes Bove, with contributions by Fabian Scheipl

### Examples

```
## get the samples
samples <- extractSamples(file.path(system.file("examples/samples", package="BayesX"),
                                         "res"))

str(samples)

## check deviance convergence
plot(samples$Deviance)

## fixed parametric effects
plot(samples$FixedEffects)

## nonparametric effects:

## handy plot function to get means and pointwise credible intervals
nonpPlot <- function(samplesMatrix,
                      ...)
{
  x <- as.numeric(colnames(samplesMatrix))

  yMeans <- colMeans(samplesMatrix)
  yCredible <- t(apply(samplesMatrix,
                      MARGIN=2,
                      FUN=quantile,
                      prob=c(0.025, 0.975),
                      na.rm=TRUE))

  matplot(x, cbind(yMeans, yCredible),
          type="l",
          lty=c(1, 2, 2),
          lwd=c(2, 1, 1),
          col=c(1, 2, 2),
          ...)
}

nonpPlot(samples$f_x1$functionSamples,
```

```
        xlab=expression(x[1]),
        ylab=expression(hat(f)(x[1])))
nonpPlot(samples$f_x2$functionSamples,
        xlab=expression(x[2]),
        ylab=expression(hat(f)(x[2])))

## spatial effect
tanzania <- read.bnd(file=system.file("examples/tanzania.bnd", package="BayesX"))
drawmap(map=tanzania,
        data=
        with(samples$f_district,
            data.frame(name=colnames(functionSamples),
                estimate=colMeans(functionSamples))),
        regionvar="name",
        plotvar="estimate")
```

---

fuse

*Combine Regions*

---

### Description

Combines a list of several regions of a map object in boundary format into a single region.

### Usage

```
fuse(map, regions, name)
```

### Arguments

map	Map object in boundary format that should be modified.
regions	Vector of regions to be combined
name	Name that should be given to the region arising from fusing the specified regions.

### Value

Map object in boundary format with the specified regions combined.

### Author(s)

Nadja Klein

### See Also

[read.bnd](#), [write.bnd](#)

## Examples

```
require("gpclib")
library("maptools")
gpclibPermit()
map <- read.bnd(system.file("examples/germany9301.bnd", package="BayesX"))
drawmap(map=map, drawnames=TRUE)

#vector of regions to be combined
regions <- c("1056","1060","1061")
#new name of combined region
newname <- "1"
newmap <- fuse(map,regions,newname)
drawmap(map=newmap,drawnames=TRUE)

#vector of regions to be combined
germany <- read.bnd(system.file("examples/germany.bnd", package="BayesX"))
drawmap(map=germany, drawnames=TRUE)
regions <- c("9371","9373","9374","9471","9472","9474","9574")
#new name of combined region
newname <- "1"
newmap <- fuse(germany,regions,newname)
drawmap(map=newmap,drawnames=TRUE)
```

---

get.centroids

*Compute Centroids of Polygons*

---

## Description

Computes all areas and centroids of the regions of a given map in boundary format.

## Usage

```
get.centroids(map)
```

## Arguments

map                    Map object in boundary format.

## Value

Matrix of area and centroids.

## Author(s)

Felix Heinzl, Thomas Kneib

**Examples**

```
germany <- read.bnd(system.file("examples/germany.bnd", package="BayesX"))
centroids <- get.centroids(germany)
centroids[1:10,]

plot(c(2100,3700),c(6800,8500),type="n", xlab="", ylab="")
for(i in 1:10){
  polygon(germany[[i]])
  region <- attr(germany,"names")[i]
  text(x=centroids[i,2]+50, y=centroids[i,3]+30, region, cex=0.7)
}
points(centroids[1:10,2:3], col='red', pch=16)
```

---

`get.neighbor`*Obtain Neighbors of Given Regions*

---

**Description**

Extracts the neighbors of a number of regions from a map in graph format.

**Usage**

```
get.neighbor(map, regions)
```

**Arguments**

<code>map</code>	Map object in graph format.
<code>regions</code>	Vector of names of regions for which the neighbors should be extracted.

**Value**

A list of vectors containing the neighbors of the elements in regions.

**Author(s)**

Felix Heinzl, Thomas Kneib

**See Also**

[add.neighbor](#), [delete.neighbor](#)

**Examples**

```
germany <- read.gra(system.file("examples/germany.gra", package="BayesX"))
get.neighbor(germany, "1001")
get.neighbor(germany, c("1001", "7339"))
```

hpd

*Computing Highest Posterior Density (HPD) Intervals***Description**

Compute approximate HPD intervals out of MCMC-samples in BayesX

**Usage**

```
hpd(data, alpha = 0.05, ...)
hpd.coda(data, alpha = 0.05, ...)
```

**Arguments**

data	Either the name of a file or a data frame containing the sample.
alpha	A numeric scalar in the interval (0,1) such that 1 - alpha is the target probability contained in the intervals.
...	Further parameters to be passed to the internal call of optim such as start and integrate such as lower.

**Details**

hpd computes the HPD interval based on a kernel density estimate of the samples. hpd.coda computes the HPD interval with the function `HPDinterval` available in package coda.

**Author(s)**

Nadja Klein

**Examples**

```
res <- read.table(system.file("examples/nonparametric_f_x_pspline_sample.raw",
                             package="BayesX"), header=TRUE)
hpd(res)
hpd.coda(res)
```

---

 Interface between nb and gra format

*Convert nb and gra format into each other*

---

**Description**

Convert neighborhood structure objects of class "nb" from R-package spdep to graph objects of class "gra" from R-package BayesX and vice versa.

**Usage**

```
nb2gra(nbObject)
gra2nb(graObject)
```

**Arguments**

```
nbObject      neighborhood structure object of class "nb"
graObject     graph object of class "gra"
```

**Value**

Equivalent object in the other format.

**Author(s)**

Daniel Sabanes Bove

**See Also**

[sp2bnd](#), [bnd2sp](#) for conversion between the geographical information formats and [read.gra](#), [write.gra](#) for the interface to the BayesX files.

**Examples**

```
## first nb to gra:
library(spdep)
library(maptools)
columbus <- readShapePoly(system.file("etc/shapes/columbus.shp", package="spdep")[1])
colNb <- poly2nb(columbus)
## ... here manual editing is possible ...
## then export to graph format
colGra <- nb2gra(colNb)

## and save in BayesX file
graFile <- tempfile()
write.gra(colGra, file=graFile)

## now back from gra to nb:
colGra <- read.gra(graFile)
newColNb <- gra2nb(colGra)
newColNb
## compare this with the original
colNb
## only the call attribute does not match (which is OK):
all.equal(newColNb, colNb,
          check.attributes=FALSE)
attr(newColNb, "call")
attr(colNb, "call")
```

---

Interface between sp and bnd format

*Convert sp and bnd format into each other*

---

### Description

Convert geographical information objects of class "SpatialPolygons" (or specializations) from R-package sp to objects of class "bnd" from R-package BayesX and vice versa.

### Usage

```
sp2bnd(spObject, regionNames, height2width, epsilon)
bnd2sp(bndObject)
```

### Arguments

spObject	object of class "SpatialPolygons" (or specializations)
regionNames	character vector of region names (parallel to the Polygons list in spObject), defaults to the IDs
height2width	ratio of total height to width, defaults to the bounding box values
epsilon	how much can two polygons differ (in maximum squared Euclidean distance) and still match each other?, defaults to machine precision
bndObject	object of class "bnd"

### Value

Equivalent object in the other format.

### Author(s)

Daniel Sabanes Bove

### See Also

[nb2gra](#), [gra2nb](#) for conversion between the neighborhood structure formats and [read.bnd](#), [write.bnd](#) for the interface to the BayesX files.

### Examples

```
## bnd to sp:
germany <- read.bnd(system.file("examples/germany2001.bnd", package="BayesX"))
spGermany <- bnd2sp(germany)

## plot the result together with the neighborhood graph
library(sp)
plot(spGermany)
library(spdep)
```



```

nbGermany <- poly2nb(spGermany)
plot(nbGermany, coords=coordinates(spGermany), add=TRUE)

## example with one region inside another
spExample <- spGermany[c("7211", "7235"), ]
plot(spExample)
plot(poly2nb(spExample), coords=coordinates(spExample), add=TRUE)

## now back from sp to bnd:
bndGermany <- sp2bnd(spGermany)
drawmap(map=bndGermany)

## compare names and number of polygons
stopifnot(identical(names(bndGermany),
                    names(germany)),
          identical(length(bndGermany),
                    length(germany)))

## compare contains-relations
surrounding <- attr(bndGermany, "surrounding")
whichInner <- which(sapply(surrounding, length) > 0L)
bndContainsData <- data.frame(inner=names(bndGermany)[whichInner],
                              outer=unlist(surrounding))

surrounding <- attr(germany, "surrounding")
whichInner <- which(sapply(surrounding, length) > 0L)
originalContainsData <- data.frame(inner=names(germany)[whichInner],
                                   outer=unlist(surrounding))

stopifnot(all(bndContainsData[order(bndContainsData$inner), ] ==
             originalContainsData[order(originalContainsData$inner), ]))

```

---

plotautocor

*Computing and Plotting Autocorrelation Functions*


---

## Description

Computes and plot autocorrelation functions for samples obtained with MCMC in BayesX

## Usage

```
plotautocor(data, ask = TRUE, lag.max=100, ...)
```

## Arguments

data	Either the name of a file or a data frame containing the sample.
ask	plotautocor will plot separate autocorrelation functions for each parameter. If ask=TRUE, the user will be prompted before showing the next plot.
lag.max	Maximum number of lags to be considered.
...	Further parameters to be passed to the internal call of plot such as ylim, etc.

**Author(s)**

Felix Heinzl, Thomas Kneib

**Examples**

```
res <- read.table(system.file("examples/nonparametric_f_x_pspline_sample.raw",
                             package="BayesX"), header=TRUE)
plotautocor(res)
plotautocor(res, lag.max=50)
```

---

plotnonp

*Plotting Nonparametric Function Estimates*

---

**Description**

Plots nonparametric function estimates obtained from BayesX

**Usage**

```
plotnonp(data, x = 2, y = c(3, 4, 5, 7, 8), ylim = NULL,
          lty = c(1, 2, 3, 2, 3), cols = rep(1, length(y)), month, year, step=12,
          xlab, ylab, ...)
```

**Arguments**

<code>data</code>	Either the name of a file or a data frame containing the estimation results.
<code>x</code>	Defines the x-axis in the plot. Either the name of a variable in <code>data</code> or the index of the corresponding column.
<code>y</code>	Defines the variables to be plotted against <code>x</code> . May be either a vector of names of variables in <code>data</code> or the corresponding indices. The default choice corresponds to the point estimate plus two confidence bands.
<code>ylim</code>	Since <code>plotnonp</code> plots multiple y-variables, it automatically determines the appropriate <code>ylim</code> to make all curves visible. Argument <code>ylim</code> allows to override this default behaviour with fixed values.
<code>lty</code>	Vector of line types used for plotting (must have the same length as <code>y</code> ). The default corresponds to solid lines for the point estimate and dashed and dotted lines for the confidence bands.
<code>cols</code>	Vector of colors used for plotting (must have the same length as <code>y</code> ). Default are black lines.
<code>month, year, step</code>	Provide specific annotation for plotting estimation results for temporal variables. <code>month</code> and <code>year</code> define the minimum time point whereas <code>step</code> specifies the type of temporal data with <code>step=4</code> , <code>step=2</code> and <code>step=1</code> corresponding to quarterly, half yearly and yearly data.
<code>xlab, ylab</code>	<code>plotnonp</code> constructs default labels that can be overwritten by these arguments
<code>...</code>	Further arguments to be passed to the interval call of <code>plot</code> such as <code>type</code> , etc.

**Author(s)**

Felix Heinzl, Andreas Brezger and Thomas Kneib

**See Also**

[drawmap](#), [plotautocor](#), [plotsample](#), [plotsurf](#)

**Examples**

```
res <- read.table(system.file("examples/nonparametric_f_x_pspline.res",
                             package="BayesX"), header=TRUE)

plotnonp(res)
plotnonp(res, x="x")
plotnonp(res, x="x", y="pmean")
plotnonp(res, x="x", y="pmed")
plotnonp(res, x="x", y="pmed", ylim=c(-2,2))
plotnonp(res, x="x", y=c("pmean", "pqu10", "pqu90"), lty=c(1,1,1),
          col=c("red", "blue", "blue"))
plotnonp(res, xlab="some variable", ylab="f(some variable)",
          main="Nonlinear effect of some variable", sub="penalised spline")

res <- read.table(system.file("examples/nonparametric2_f_time_pspline.res",
                             package="BayesX"), header=TRUE)

plotnonp(res)
plotnonp(res, month=1, year=1980, step=12)

res <- res[1:18,]
plotnonp(res, month=1, year=1980, step=12)
```

---

plotsample

*Plotting Sampling Paths*


---

**Description**

Plots sampling paths obtained with MCMC-sampling in BayesX

**Usage**

```
plotsample(data, ask = TRUE, ...)
plotsample.coda(data, ask = TRUE, ...)
```

**Arguments**

data	Either the name of a file or a data frame containing the sample.
ask	plotsample will plot separate sampling paths for each parameter. If ask=TRUE, the user will be prompted before showing the next plot.
...	Further parameters to be passed to the internal call of plot such as ylim, etc.

**Details**

plotsample simply plots sampling paths while plotsampe . coda makes use of the plotting facilities available in package coda.

**Author(s)**

Felix Heinzl, Andreas Brezger, Thomas Kneib

**See Also**

[drawmap](#), [plotautocor](#), [plotnonp](#), [plotsurf](#),

**Examples**

```
res <- read.table(system.file("examples/nonparametric_f_x_pspline_sample.raw",
                             package="BayesX"), header=TRUE)
plotsample(res)
```

---

plotsurf

*Visualise Surface Estimates*


---

**Description**

Visualises surface estimates obtained with BayesX.

**Usage**

```
plotsurf(data, x=2, y=3, z=4, mode=1, ticktype="detailed",
          expand=0.75, d=100, theta=-30, phi=25, ...)
akimaPermitStatus()
akimaPermit()
```

**Arguments**

data	Either the name of a file or a data frame containing the estimation results.
x	Defines the x-axis in the plot. Either the name of a variable in data or the index of the corresponding column.
y	Defines the y-axis in the plot. Either the name of a variable in data or the index of the corresponding column.
z	Defines the z-axis in the plot. Either the name of a variable in data or the index of the corresponding column.
mode	plotsurf is mostly a convenient interface to the functions persp (mode=1), image (mode=2) and contour (mode=3).
ticktype, expand, d, theta, phi	Overwrite the default behaviour of persp
...	Further parameters that are parsed to the internal call to persp, image or contour

**Details**

The `akimaPermit` function is used to choose to permit the use of `akima` if installed, and `akimaPermitStatus` reports its status. The licence for `akima` is not Free or Open Source and explicitly forbids commercial use. See the `akima` licence file for details.

**Author(s)**

Felix Heinzl, Thomas Kneib

**See Also**

[drawmap](#), [plotautocor](#), [plotsample](#), [plotnonp](#)

**Examples**

```
res <- read.table(system.file("examples/surface_f_x1_x2_pspline.res",
                             package="BayesX"), header=TRUE)

# call akimaPermit to allow use of akima despite its licence that restricts
# usage to non-commercial purposes
akimaPermit()
plotsurf(res)
plotsurf(res, mode=2)
plotsurf(res, mode=3)

plotsurf(res, x="x1", y="x2", z="pmed")

plotsurf(res, ticktype="simple")

plotsurf(res, main="3D-Plot", xlab="myx", ylab="myy", zlab="f(myx,myy)")
```

---

read.bnd

*Read Geographical Information in Boundary Format*

---

**Description**

Reads the geographical information provided in a file in boundary format (see Ch. 5 of the BayesX Reference Manual) and stores it in a map object.

**Usage**

```
read.bnd(file, sorted=FALSE)
```

**Arguments**

<code>file</code>	Name of the boundary file to be read.
<code>sorted</code>	Should the regions be ordered by the numbers specifying the region names ( <code>sorted=TRUE</code> )?

**Value**

Returns a list of polygons that form the map. Additional attributes are

surrounding	Parallel list where for each polygon, the name of a possible surrounding region is saved.
height2width	Ratio between height and width of the map. Allows customised drawing and storage in files by specifying the appropriate height and width.
class	Indicates whether the map is stored in boundary format (bnd) or graph format (gra). Maps returned by <code>read.bnd</code> are of class <code>bnd</code>

**Author(s)**

Daniel Sabanes Bove, Felix Heinzl, Thomas Kneib, Andreas Brezger

**References**

BayesX Reference Manual. Available from <http://www.bayesx.org/>

**See Also**

[write.bnd](#), [drawmap](#), [read.gra](#), [write.gra](#)

**Examples**

```
germany <- read.bnd(system.file("examples/germany.bnd", package="BayesX"))
drawmap(map=germany)
attributes(germany)
```

```
germany <- read.bnd(system.file("examples/germany2001.bnd", package="BayesX"))
drawmap(map=germany)
attributes(germany)
```

---

read.gra

*Read Geographical Information in Graph Format*

---

**Description**

Reads the geographical information provided in a file in graph format (see Ch. 5 of the BayesX Reference Manual) and stores it in a map object.

**Usage**

```
read.gra(file, sorted=FALSE)
```

**Arguments**

file	Name of the graph file to be read.
sorted	Should the regions be ordered by the numbers specifying the region names ( <code>sorted=TRUE</code> )?

**Value**

Returns an adjacency matrix that represents the neighborhood structure defined in the graph file. Additional attributes are

<code>dim</code>	Dimension of the (square) adjacency matrix.
<code>dimnames</code>	List of region names corresponding to rows and columns of the adjacency matrix.
<code>class</code>	Indicates whether the map is stored in boundary format ( <code>bnd</code> ) or graph format ( <code>gra</code> ). Maps returned by <code>read.gra</code> are of class <code>gra</code>

**Author(s)**

Thomas Kneib, Felix Heinzl

**References**

BayesX Reference Manual. Available from <http://www.bayesx.org/>

**See Also**

[write.gra](#), [read.bnd](#), [write.bnd](#), [get.neighbor](#), [add.neighbor](#), [delete.neighbor](#)

**Examples**

```
germany <- read.gra(system.file("examples/germany.gra", package="BayesX"))
attributes(germany)
```

---

shp2bnd

*convert a shape-file into a boundary object*

---

**Description**

Converts the geographical information provided in a shape-file into a boundary object (see Ch. 5 of the Reference Manual)

**Usage**

```
shp2bnd(shpname, regionnames, check.is.in = TRUE)
```

**Arguments**

<code>shpname</code>	Base filename of the shape-file (including path)
<code>regionnames</code>	Either a vector of region names or the name of the variable in the dbf-file representing these names
<code>check.is.in</code>	Test whether some regions are surrounded by other regions ( <code>FALSE</code> speeds up the execution time but may result in a corrupted bnd-file)

**Value**

Returns a boundary object, i.e. a list of polygons that form the map. See [read.bnd](#) for more information on the format.

**Author(s)**

Felix Heinzl, Daniel Sabanes Bove, Thomas Kneib with contributions by Michael Hoehle and Frank Sagerer

**References**

BayesX Reference Manual. Available from <http://www.bayesx.org>

**See Also**

[write.bnd](#), [drawmap](#), [read.bnd](#)

**Examples**

```
## read shapefile into bnd object
shpName <- sub(pattern="(.*)\.dbf", replacement="\1",
              x=system.file("examples/northamerica_adm0.dbf",
                           package="BayesX"))
north <- shp2bnd(shpname=shpName, regionnames="COUNTRY")

## draw the map
drawmap(map=north)

## compare with shipped bnd file
shippedBnd <- read.bnd(system.file("examples/northamerica.bnd", package="BayesX"))
stopifnot(all.equal(north, shippedBnd))
```

---

smooth.bnd

*Round Boundary Information*

---

**Description**

Rounds the boundary information in a map object in boundary format to a specified precision.

**Usage**

```
smooth.bnd(map, digits = 2, scale = 1)
```

**Arguments**

map	Map object in boundary format that should be modified.
digits	Number of digits to round to.
scale	Scaling factor that should be applied for rounding. For example, with scale=0.1 all polygons are divided by 10 before rounding.



**Value**

Map object in boundary format rounded to the specified precision.

**Author(s)**

Felix Heinzl, Thomas Kneib

**See Also**

[read.bnd](#), [write.bnd](#)

---

write.bnd

*Saving Maps in Boundary Format*

---

**Description**

Writes the information of a map object to a file (in boundary format)

**Usage**

```
write.bnd(map, file, replace = FALSE)
```

**Arguments**

map	Map object to be saved (should be in boundary format).
file	Name of the file to write to
replace	Should an existing file be overwritten with the new version?

**Author(s)**

Thomas Kneib, Felix Heinzl

**References**

BayesX Reference Manual. Available from <http://www.bayesx.org>

**See Also**

[write.gra](#), [read.gra](#), [read.bnd](#)

---

`write.gra`*Saving Maps in Graph Format*

---

**Description**

Writes the information of a map object to a file (in graph format)

**Usage**

```
write.gra(map, file, replace = FALSE)
```

**Arguments**

<code>map</code>	Map object to be saved (should be in graph format, see <a href="#">bnd2gra</a> for the conversion of boundary format to graph format).
<code>file</code>	Name of the file to write to
<code>replace</code>	Should an existing file be overwritten with the new version?

**Author(s)**

Thomas Kneib, Felix Heinzl

**References**

BayesX Reference Manual. Available from <http://www.bayesx.org>

**See Also**

[write.bnd](#), [read.gra](#), [read.bnd](#)

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