

# Package ‘RBaM’

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**Type** Package

**Title** Bayesian Modeling: Estimate a Computer Model and Make Uncertain Predictions

**Version** 1.0.1

**Description** An interface to the 'BaM' (Bayesian Modeling) engine, a 'Fortran'-based executable aimed at estimating a model with a Bayesian approach and using it for prediction, with a particular focus on uncertainty quantification. Classes are defined for the various building blocks of 'BaM' inference (model, data, error models, Markov Chain Monte Carlo (MCMC) samplers, predictions).

The typical usage is as follows:

- (1) specify the model to be estimated;
- (2) specify the inference setting (dataset, parameters, error models...);
- (3) perform Bayesian-MCMC inference;
- (4) read, analyse and use MCMC samples;
- (5) perform prediction experiments.

Technical details are available (in French) in

Renard (2017) <<https://hal.science/hal-02606929v1>>.

Examples of applications include

Mansanarez et al. (2019) <[doi:10.1029/2018WR023389](https://doi.org/10.1029/2018WR023389)>,

Le Coz et al. (2021) <[doi:10.1002/hyp.14169](https://doi.org/10.1002/hyp.14169)>,

Perret et al. (2021) <[doi:10.1029/2020WR027745](https://doi.org/10.1029/2020WR027745)>,

Darienzo et al. (2021) <[doi:10.1029/2020WR028607](https://doi.org/10.1029/2020WR028607)> and

Perret et al. (2023) <[doi:10.1061/JHEND8.HYENG-13101](https://doi.org/10.1061/JHEND8.HYENG-13101)>.

**License** GPL-3

**Encoding** UTF-8

**LazyData** true

**URL** <https://github.com/BaM-tools/RBaM>

**BugReports** <https://github.com/BaM-tools/RBaM/issues>

**Depends** R (>= 4.0.0)

**Imports** ggplot2, gridExtra, R.utils, utils, grDevices, stats, rjson, rlang, tools, tidyverse

**Suggests** rmarkdown

**RoxygenNote** 7.2.3

**NeedsCompilation** no

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**BaM***Run BaM***Description**

Run BaM.exe

**Usage**

```
BaM(
  workspace,
  mod,
  data,
  remnant = rep(list(remnantModelError()), mod$nY),
  mcmc = mcmcOptions(),
  cook = mcmcCooking(),
  summary = mcmcSummary(),
  residuals = residualOptions(),
  pred = NULL,
  doCalib = TRUE,
  doPred = FALSE,
  na.value = -9999,
  run = TRUE,
  preClean = FALSE,
  dir.exe = .BAM_PATH,
  name.exe = "BaM",
  predMaster_fname = "Config_Pred_Master.txt"
)
```

**Arguments**

workspace	Character, directory where config and result files are stored.
mod	model object, the model to be calibrated
data	dataset object, calibration data
remnant	list of remnantModelError objects. WARNING: make sure you use a list of length mod\$nY (even if mod\$nY=1!)
mcmc	mcmcOptions object, MCMC simulation number and options
cook	mcmcCooking object, properties of MCMC cooking (burn and slice)
summary	mcmcSummary object, properties of MCMC summary

residuals	residualOptions object, properties of residual analysis
pred	list of prediction objects, properties of prediction experiments
doCalib	Logical, do Calibration? (mcmc+cooking+summary+residuals)
doPred	Logical, do Prediction?
na.value	numeric, value used for NAs when writing the dataset in BaM format
run	Logical, run BaM? if FALSE, just write config files.
preClean	Logical, start by cleaning up workspace? Be careful, this will delete all files in the workspace, including old results!
dir.exe	Character, directory where BaM executable stands.
name.exe	Character, name of the executable without extension ('BaM' by default).
predMaster_fname	Character, name of configuration file pointing to all prediction experiments.

### Value

Nothing: just write config files and runs the executable.

### Examples

```
# Fitting a rating curve - see https://github.com/BaM-tools/RBaM
workspace=tempdir()
D=dataset(X=SauzeGaugings['H'],Y=SauzeGaugings['Q'],Yu=SauzeGaugings['uQ'],data.dir=workspace)
# Parameters of the low flow section control: activation stage k, coefficient a and exponent c
k1=parameter(name='k1',init=-0.5,prior.dist='Uniform',prior.par=c(-1.5,0))
a1=parameter(name='a1',init=50,prior.dist='LogNormal',prior.par=c(log(50),1))
c1=parameter(name='c1',init=1.5,prior.dist='Gaussian',prior.par=c(1.5,0.05))
# Parameters of the high flow channel control: activation stage k, coefficient a and exponent c
k2=parameter(name='k2',init=1,prior.dist='Gaussian',prior.par=c(1,1))
a2=parameter(name='a2',init=100,prior.dist='LogNormal',prior.par=c(log(100),1))
c2=parameter(name='c2',init=1.67,prior.dist='Gaussian',prior.par=c(1.67,0.05))
# Define control matrix: columns are controls, rows are stage ranges.
controlMatrix=rbind(c(1,0),c(0,1))
# Stitch it all together into a model object
M=model(ID='BaRatin',
         nX=1,nY=1, # number of input/output variables
         par=list(k1,a1,c1,k2,a2,c2), # list of model parameters
         xtra=xtraModelInfo(object=controlMatrix)) # use xtraModelInfo() to pass the control matrix
# Call BaM to write configuration files. To actually run BaM, use run=TRUE,
# but BaM executable needs to be downloaded first (use downloadBaM())
BaM(workspace=workspace,mod=M,data=D,run=FALSE)
```

---

dataset	<i>dataset object constructor.</i>
---------	------------------------------------

---

## Description

Creates a new instance of a 'dataset' object

## Usage

```
dataset(  
  X,  
  Y,  
  data.dir = getwd(),  
  data.fname = "CalibrationData.txt",  
  fname = "Config_Data.txt",  
  Xu = NULL,  
  Xb = NULL,  
  Xb.indx = NULL,  
  Yu = NULL,  
  Yb = NULL,  
  Yb.indx = NULL,  
  VAR.indx = NULL  
)
```

## Arguments

X	data frame, observed input variables.
Y	data frame, observed output variables (same number of rows as X).
data.dir	Character, directory where a copy of the dataset will be written if required. Default is the current working directory, but you may prefer to use the BaM workspace.
data.fname	Character, data file name.
fname	Character, configuration file name.
Xu	data frame, random uncertainty in X, expressed as a standard deviation. Same dimension as X.
Xb	data frame, systematic uncertainty in X, expressed as a standard deviation. Same dimension as X.
Xb.indx	data frame, index of systematic errors in X. Same dimension as X.
Yu	data frame, random uncertainty in Y, expressed as a standard deviation. Same dimension as Y.
Yb	data frame, systematic uncertainty in Y, expressed as a standard deviation. Same dimension as Y.
Yb.indx	data frame, index of systematic errors in Y. Same dimension as Y.
VAR.indx	data frame, indices used for defining how VAR parameters vary.

**Value**

An object of class 'dataset'.

**Examples**

```
X=data.frame(input1=rnorm(100),input2=rnorm(100))
Y=data.frame(output=X$input1+0.8*X$input2+0.1*rnorm(100))
workspace=tempdir()
d <- dataset(X=X,Y=Y,data.dir=workspace)
```

densityPlot

*densityPlot***Description**

returns a histogram+density ggplot (or a list thereof if several columns in sim)

**Usage**

```
densityPlot(sim, xlab = "values", col = "black")
```

**Arguments**

sim	vector or matrix or data frame, MCMC simulations
xlab	Character, label of x-axis to be used if sim has no names
col	Color

**Value**

A ggplot (or a list thereof if several columns in sim)

**Examples**

```
# Create Monte Carlo samples
n=1000
sim=data.frame(p1=rnorm(n),p2=rlnorm(n),p3=runif(n))
# create density plot for each component
figures=densityPlot(sim)
```

---

downloadBaM*BaM downloader*

---

**Description**

Download BaM executable

**Usage**

```
downloadBaM(  
  destFolder,  
  url = NULL,  
  os = Sys.info()["sysname"],  
  quiet = FALSE,  
  ...  
)
```

**Arguments**

destFolder	character string, folder where BaM executable will be downloaded.
url	character string, the url from which BaM should be downloaded. When NULL, the url is determined automatically by using GitHub API to determine the latest release and the file corresponding to the OS.
os	character string, operating system, e.g. 'Linux', 'Windows' or 'Darwin'.
quiet	logical, if TRUE, suppress status messages.
...	arguments passed to function 'download.file'

**Value**

nothing - just download the file.

**Examples**

```
try(downloadBaM(destFolder=tempdir()))
```

---

getAwPfromBathy*Bathymetry interpreter*

---

**Description**

Compute Area A(h), width w(h) and wet perimeter P(h) from a bathymetry profile (a,z).

**Usage**

```
getAwPfromBathy(
  bathy,
  hgrid = seq(min(bathy[, 2]), max(bathy[, 2]), diff(range(bathy[, 2]))/1000),
  segmentLength = sum(sqrt(apply(apply(bathy, 2, diff)^2, 1, sum)))/1000
)
```

**Arguments**

**bathy** data frame, 2 columns containing abscissa (increasing values) and stage.  
**hgrid** numeric vector, grid of h values where A, w and P are computed. By default 1000 values in the range of bathymetry's z.  
**segmentLength** numeric, segment length for bathymetry subsampling. By default 1/1000 of the total bathymetry's perimeter.

**Value**

A 4-column dataframe containing h, A(h), w(h) and P(h)

**Examples**

```
bathy=data.frame(a=c(0,0,0,1,2,2,4,6,8),h=c(3,2,0,-0.5,0,2,2.0001,2.3,3))
plot(bathy,type='l')
df=getAwPfromBathy(bathy)
plot(df$h,df$A,type='l')
plot(df$h,df$w,type='l')
plot(df$h,df$P,type='l')
```

getCatalogue

*BaM catalogue*

**Description**

Distributions and models available in BaM

**Usage**

```
getCatalogue(printOnly = FALSE)
```

**Arguments**

**printOnly** Logical, should the catalogue be returned or only printed?

**Value**

If `printOnly==FALSE`, a list with the following fields:

**distributions** available univariate distributions.

**models** available models.

**Examples**

```
catalogue <- getCatalogue()
getCatalogue(printOnly=TRUE)
```

---

getNames

*Get object names*

---

**Description**

getNames from an object or a list of objects having a \$name field (e.g. parameters)

**Usage**

```
getNames(loo, name = "name")
```

**Arguments**

loo	List Of Objects
name	character, string denoting the name field

**Value**

A character vector containing names

**Examples**

```
pars <- list(parameter(name='par1',init=0),
              parameter(name='par2',init=0),
              parameter(name='Third parameter',init=0))
getNames(pars)
```

---

getParNames

*Get parameter names*

---

**Description**

Get parameter names for a distribution d

**Usage**

```
getParNames(d)
```

**Arguments**

d	Character (possibly vector), distribution (possibly distributions)
---	--

**Value**

A character vector with parameter names.

**Examples**

```
parnames <- getParNames('GEV')
npar <- length(getParNames('Gumbel'))
```

---

**mcmcCooking**

*mcmcCooking constructor.*

---

**Description**

Creates a new instance of a 'mcmcCooking' object

**Usage**

```
mcmcCooking(
  fname = "Config_Cooking.txt",
  result.fname = "Results_Cooking.txt",
  burn = 0.5,
  nSlim = 10
)
```

**Arguments**

<code>fname</code>	Character, configuration file name.
<code>result.fname</code>	Character, result file name.
<code>burn</code>	numeric, burn factor, $\geq 0$ and $< 1$ . 0.4 means the first 40 percent of MCMC samples are discarded).
<code>nSlim</code>	Integer, slimming period: 10 means only one MCMC sample every 10 is kept (after burning).

**Value**

An object of class 'mcmcCooking'.

**Examples**

```
m <- mcmcCooking()
```

---

<code>mcmcOptions</code>	<i>mcmcOptions object constructor.</i>
--------------------------	--

---

## Description

Creates a new instance of a 'mcmcOptions' object

## Usage

```
mcmcOptions(
  fname = "Config_MCMC.txt",
  result.fname = "Results_MCMC.txt",
  nAdapt = 100,
  nCycles = 100,
  minMoveRate = 0.1,
  maxMoveRate = 0.5,
  downMult = 0.9,
  upMult = 1.1,
  multFactor = 0.1,
  manualMode = FALSE,
  thetaStd = 9999,
  gammaStd = list(9999)
)
```

## Arguments

<code>fname</code>	Character, configuration file name.
<code>result.fname</code>	Character, result file name.
<code>nAdapt</code>	Integer, adaptation period: jump sizes are increased/decreased every Nadapt iterations to comply with the desired moving rates.
<code>nCycles</code>	Integer, number of adaptation cycles (total number of iterations is hence Nadapt * Ncycles).
<code>minMoveRate</code>	Numeric in (0;1), lower bound for the desired move rate interval.
<code>maxMoveRate</code>	Numeric in (0;1), upper bound for the desired move rate interval.
<code>downMult</code>	Numeric in (0:1), multiplication factor used to decrease jump size when move rate is too low.
<code>upMult</code>	Numeric (>1, avoid 1/dowMult) multiplication factor used to increase jump size when move rate is too high.
<code>multFactor</code>	Numeric >0, multiplicative factor to set initial jump standard deviations to multFactor*initValue1 (AUTO mode).
<code>manualMode</code>	logical, should jump standard deviations be entered manually?
<code>thetaStd</code>	Numeric vector (>0), jump standard deviations for model parameters theta (MANUAL mode).
<code>gammaStd</code>	list of numeric vectors (>0), size = number of output variables of the model. Jump standard deviations for structural error parameters gamma of each output variable (MANUAL mode).

**Value**

An object of class 'mcmcOptions'.

**Examples**

```
m <- mcmcOptions()
```

---

**mcmcSummary**

*mcmcSummary constructor.*

---

**Description**

Creates a new instance of a 'mcmcSummary' object

**Usage**

```
mcmcSummary(
  fname = "Config_Summary.txt",
  result.fname = "Results_Summary.txt",
  DIC.fname = "Results_DIC.txt",
  xtendedMCMC.fname = ""
)
```

**Arguments**

<code>fname</code>	Character, configuration file name.
<code>result.fname</code>	Character, summary file name.
<code>DIC.fname</code>	Character, DIC file name. Not computed if empty string.
<code>xtendedMCMC.fname</code>	Character, xtended MCMC file name. Not written if empty string.

**Value**

An object of class 'mcmcSummary'.

**Examples**

```
m <- mcmcSummary()
```

---

MeyrasGaugingsMeyras Gaugings

---

## Description

Stage-discharge gaugings from the hydrometric station 'the Ardèche River at Meyras'. See <https://en.wikipedia.org/wiki/Ardèche> for a description of the river See <https://doi.org/10.1029/2018WR023389> for an article using this dataset

## Usage

MeyrasGaugings

## Format

A data frame with 104 rows and 4 variables:

**h** Stage (m)

**Q** Discharge (m<sup>3</sup>/s)

**uQ** Discharge uncertainty (m<sup>3</sup>/s) expressed as a standard deviation

**Period** Stability period on which a single rating curve can be used

---

modelmodel object constructor.

---

## Description

Creates a new instance of a 'model' object

## Usage

```
model(
  fname = "Config_Model.txt",
  ID = "Linear",
  nX = 1,
  nY = 1,
  par = list(parameter("Xeffect", 1, prior.dist = "FlatPrior")),
  xtra = xtraModelInfo()
)
```

**Arguments**

<code>fname</code>	Character, configuration file name.
<code>ID</code>	Character, model ID. Type <code>'getCatalogue()'</code> for available models.
<code>nX</code>	Integer, number of input variables.
<code>nY</code>	Integer, number of output variables.
<code>par</code>	list of parameter objects, parameters of the model.
<code>xtra</code>	xtraModelInfo object.

**Value**

An object of class `'model'`.

**Examples**

```
# default linear regression model Y=aX+b
mod <- model()
# BaRatin model for a single-control rating curve Y=a(X-b)^c
mod <- model(ID='BaRatin',nX=1,nY=1,
               par=list(parameter('a',10,prior.dist='LogNormal',prior.par=c(log(10),0.1)),
                        parameter('b',-1,prior.dist='Gaussian',prior.par=c(-1,1)),
                        parameter('c',5/3,prior.dist='Gaussian',prior.par=c(5/3,0.05))),
               xtra=xtraModelInfo(object=matrix(1,nrow=1,ncol=1)))
```

<code>parameter</code>	<i>parameter object constructor.</i>
------------------------	--------------------------------------

**Description**

Creates a new instance of a `'parameter'` object

**Usage**

```
parameter(name, init, prior.dist = "FlatPrior", prior.par = NULL)
```

**Arguments**

<code>name</code>	character, parameter name.
<code>init</code>	numeric, initial guess.
<code>prior.dist</code>	character, prior distribution.
<code>prior.par</code>	numeric vector, prior parameters

**Value**

An object of class `'parameter'`.

**Examples**

```
p <- parameter(name='par',init=0,prior.dist='Gaussian',prior.par=c(0,1))
```

---

<code>parameter_VAR</code>	<i>Varying parameter object constructor.</i>
----------------------------	--

---

## Description

Creates a new instance of a 'parameter\_VAR' object

## Usage

```
parameter_VAR(
  name,
  index,
  d,
  init,
  prior.dist = rep("FlatPrior", length(init)),
  prior.par = rep(list(NULL), length(init))
)
```

## Arguments

<code>name</code>	character, parameter name.
<code>index</code>	character, name of column in VAR.idx (see ?dataset) containing the index for this varying parameter
<code>d</code>	dataset object, the dataset containing (amongst other things) the index above
<code>init</code>	numeric vector, initial guesses for each instance of the VAR parameter.
<code>prior.dist</code>	character vector, prior distribution for each instance of the VAR parameter.
<code>prior.par</code>	list of numeric vectors, prior parameters for each instance of the VAR parameter

## Value

An object of class 'parameter\_VAR'.

## Examples

```
X=data.frame(input1=rnorm(100),input2=rnorm(100))
Y=data.frame(output=X$input1+0.8*X$input2+0.1*rnorm(100))
VAR.idx=data.frame(idx=c(rep(1,50),rep(2,50)))
workspace=tempdir()
d <- dataset(X=X,Y=Y,data.dir=workspace,VAR.idx=VAR.idx)
p <- parameter_VAR(name='par',index='idx',d=d,
                     init=c(-1,1,2),
                     prior.dist=c('Gaussian','FlatPrior','Triangle'),
                     prior.par=list(c(-1,1),NULL,c(2,0,5)))
```

prediction	<i>prediction object constructor.</i>
------------	---------------------------------------

## Description

Creates a new instance of a 'prediction' object

## Usage

```
prediction(
  X,
  spagFiles,
  data.dir = getwd(),
  data.fnames = paste0("X", 1:length(X), ".pred"),
  fname = paste0("Config_Pred_", paste0(sample(c(letters, LETTERS, 0:9), 6), collapse =
    ""), ".txt"),
  doParametric = FALSE,
  doStructural = rep(FALSE, length(spagFiles)),
  transposeSpag = TRUE,
  priorNsim = NULL,
  envFiles = paste0(tools::file_path_sans_ext(spagFiles), ".env"),
  consoleProgress = TRUE,
  spagFiles_state = NULL,
  transposeSpag_state = TRUE,
  envFiles_state = switch(is.null(spagFiles_state) + 1,
    paste0(tools::file_path_sans_ext(spagFiles_state), ".env"), NULL),
  parSamples = NULL
)
```

## Arguments

X	data frame or list of dataframes / matrices, representing the values taken by the input variables. <ul style="list-style-type: none"> <li>• If X is a dataframe, then each column is interpreted as one input variable, and consequently inputs are not replicated (=&gt; no input uncertainty).</li> <li>• If X is a list, then each element of the list is a matrix associated with one input variable, and the columns of this matrix are replications (=&gt; input uncertainty is propagated). All matrices in the list should have the same number of rows, columns are recycled if needed.</li> </ul>
spagFiles	Character vector (size nY, the number of output variables). Name of the files containing the spaghetti for each output variable. NOTE: provide file names only, not full paths. Using a '.spag' extension is a good practice.
data.dir	Character, directory where a copies of the dataset X will be written if required (1 file per variable in X). Default is the current working directory, but you may prefer to use the BaM workspace.
data.fnames	Character, data file names.

fname	Character, configuration file name.
doParametric	Logical, propagate parametric uncertainty? If FALSE, maxpost parameters are used.
doStructural	Logical, propagate structural uncertainty for each output variable? (size nY)
transposeSpag	Logical. If FALSE, spaghettis are written horizontally (row-wise), otherwise they will be transposed so that each spaghetti is a column.
priorNsim	Integer, number of samples from the prior distribution for 'prior prediction' experiments. If negative or NULL (default), posterior samples are used.
envFiles	Character vector (size nY, the number of output variables). Name of the files containing the envelops (e.g. prediction intervals) computed from the spaghettis for each output variable. By default, same name as spaghetti files but with a '.env' extension. If NULL, envelops are not computed.
consoleProgress	Logical, print progress in BaM.exe console?
spagFiles_state	Character vector (size nState, the number of state variables), same as spagFiles but for states rather than outputs. If NULL, states are not predicted. Note that only parametric uncertainty is propagated for state variables since they are not observed. Consequently, structural uncertainty = 0 and total uncertainty = parametric uncertainty.
transposeSpag_state	Logical. Same as transposeSpag, but for states rather than outputs.
envFiles_state	Character vector (size nState, the number of state variables), same as envFiles, but for states rather than outputs.
parSamples	data frame, parameter samples that will replace the MCMC-generated one for this prediction.

### Value

An object of class 'prediction'.

### Examples

```
#-----
# Example using the twoPopulations dataset, containing 101 values for
# 3 input variables (time t, temperature at site 1 T1, temperature at site 2 T2)
# and 2 output variables (population at site 1 P1, population at site 2 P2).
pred=prediction(X=twoPopulations[,1:3],spagFiles=c('P1.spag','P2.spag'))
#-----
# Alternative example showing how to propagate uncertainty in some of
# the input variables (here, temperatures T1 and T2)
# Create 100 noisy replicates for T1, representing uncertainty
T1rep=matrix(rnorm(101*100,mean=twoPopulations$T1, sd=0.1),nrow=101,ncol=100)
# Same for T2
T2rep=matrix(rnorm(101*100,mean=twoPopulations$T2, sd=0.1),nrow=101,ncol=100)
# Create prediction object
pred=prediction(X=list(twoPopulations$t,T1rep,T2rep),spagFiles=c('P1.spag','P2.spag'))
```

---

readMCMC*MCMC Reader*

---

**Description**

Read raw MCMC samples, return cooked (burnt & sliced) ones

**Usage**

```
readMCMC(
  file = "Results_Cooking.txt",
  burnFactor = 0,
  slimFactor = 1,
  sep = "",
  reportFile = NULL,
  panelPerCol = 10,
  panelHeight = 3,
  panelWidth = 23/panelPerCol
)
```

**Arguments**

file	Character, full path to MCMC file.
burnFactor	Numeric, burn factor. 0.1 means the first 10 are discarded.
slimFactor	Integer, slim factor. 10 means that only one iteration every 10 is kept.
sep	Character, separator used in MCMC file.
reportFile	Character, full path to pdf report file, not created if NULL
panelPerCol	Integer, max number of panels per column
panelHeight	Numeric, height of each panel
panelWidth	Numeric, width of each panel

**Value**

A data frame containing the cooked mcmc samples.

**Examples**

```
# Create Monte Carlo samples and write them to file
n=4000
sim=data.frame(p1=rnorm(n),p2=rlnorm(n),p3=runif(n))
workspace=tempdir()
write.table(sim,file=file.path(workspace,'MCMC.txt'),row.names=FALSE)
# Read file, burn the first half and keep every other row
M=readMCMC(file=file.path(workspace,'MCMC.txt'),burnFactor=0.5,slimFactor=2)
dim(M)
```

---

`remnantErrorModel`      *remnantErrorModel object constructor.*

---

**Description**

Creates a new instance of a 'remnantErrorModel' object

**Usage**

```
remnantErrorModel(
  fname = "Config_RemnantSigma.txt",
  funk = "Linear",
  par = list(parameter("g1", 1, prior.dist = "FlatPrior+"), parameter("g2", 0.1,
    prior.dist = "FlatPrior+"))
)
```

**Arguments**

<code>fname</code>	Character, configuration file name.
<code>funk</code>	Character, function f used in remnant sdev = f(Ysim). Available: 'Constant', 'Proportional', 'Linear' (default), 'Exponential', 'Gaussian'.
<code>par</code>	list of parameter objects, parameters of the function above. respectively, npar=1,1,2,3,3

**Value**

An object of class 'remnantErrorModel'.

**Examples**

```
r <- remnantErrorModel()
```

---

`residualOptions`      *residualOptions constructor.*

---

**Description**

Creates a new instance of a 'residualOptions' object

**Usage**

```
residualOptions(
  fname = "Config_Residuals.txt",
  result.fname = "Results_Residuals.txt"
)
```

**Arguments**

<code>fname</code>	Character, configuration file name.
<code>result.fname</code>	Character, result file name.

**Value**

An object of class 'residualOptions'.

**Examples**

```
r <- residualOptions()
```

<code>runOptions</code>	<i>runOptions constructor.</i>
-------------------------	--------------------------------

**Description**

Creates a new instance of a 'runOptions' object

**Usage**

```
runOptions(
  fname = "Config_RunOptions.txt",
  doMCMC = TRUE,
  doSummary = TRUE,
  doResiduals = TRUE,
  doPrediction = FALSE
)
```

**Arguments**

<code>fname</code>	Character, configuration file name.
<code>doMCMC</code>	logical, do MCMC sampling?
<code>doSummary</code>	logical, do MCMC summarizing?
<code>doResiduals</code>	logical, do residuals analysis?
<code>doPrediction</code>	logical, do prediction experiments?

**Value**

An object of class 'runOptions'.

**Examples**

```
o <- runOptions()
```

---

SauzeGaugings

*Sauze Gaugings*

---

### Description

Stage-discharge gaugings from the hydrometric station 'the Ardèche River at Sauze-St-Martin'. See [https://en.wikipedia.org/wiki/Ardèche\\_\(river\)](https://en.wikipedia.org/wiki/Ardèche_(river)) for a description of the river See <https://hal.science/hal-00934237> for an article using this dataset

### Usage

SauzeGaugings

### Format

A data frame with 38 rows and 3 variables:

**H** Stage (m)

**Q** Discharge (m<sup>3</sup>/s)

**uQ** Discharge uncertainty (m<sup>3</sup>/s) expressed as a standard deviation

---

setPathToBaM

*Path to BaM*

---

### Description

Set path to BaM executable

### Usage

```
setPathToBaM(dir.exe, quiet = FALSE)
```

### Arguments

**dir.exe** character string, folder where BaM executable is located. A NULL value resets BaM directory to 'unknown' by removing the config folder where it was stored on your computer (use `'tools::R_user_dir(package="RBaM",which="config")'` to locate this folder).

**quiet** logical, if TRUE, suppress status messages.

### Value

nothing - just write a config file.

### Examples

```
setPathToBaM(dir.exe=tempdir())
```

**toString.dataset**      *dataset to string*

### Description

Convert an object of class 'dataset' into a ready-to-write vector of string

### Usage

```
## S3 method for class 'dataset'
toString(x, ...)
```

### Arguments

x	dataset object, object to be converted.
...	Optional arguments.

### Value

A string ready to be printed or written.

### Examples

```
X=data.frame(input1=rnorm(100),input2=rnorm(100))
Y=data.frame(output=X$input1+0.8*X$input2+0.1*rnorm(100))
workspace=tempdir()
d <- dataset(X=X,Y=Y,data.dir=workspace)
toString(d)
```

**toString.mcmcCooking**    *mcmcCooking to string*

### Description

Convert an object of class 'mcmcCooking' into a ready-to-write vector of string

### Usage

```
## S3 method for class 'mcmcCooking'
toString(x, ...)
```

### Arguments

x	mcmcCooking object, object to be converted.
...	Optional arguments.

**Value**

A string ready to be printed or written.

**Examples**

```
toString(mcmcCooking())
```

---

toString.mcmcOptions    *mcmcOptions to string*

---

**Description**

Convert an object of class 'mcmcOptions' into a ready-to-write vector of string

**Usage**

```
## S3 method for class 'mcmcOptions'  
toString(x, ...)
```

**Arguments**

x                mcmcOptions object, object to be converted.  
...                Optional arguments.

**Value**

A string ready to be printed or written.

**Examples**

```
toString(mcmcOptions())
```

---

toString.mcmcSummary    *mcmcSummary to string*

---

**Description**

Convert an object of class 'mcmcSummary' into a ready-to-write vector of string

**Usage**

```
## S3 method for class 'mcmcSummary'  
toString(x, ...)
```

**Arguments**

- x mcmcSummary object, object to be converted.
- ... Optional arguments.

**Value**

A string ready to be printed or written.

**Examples**

```
toString(mcmcSummary())
```

---

toString.model      *model to string*

---

**Description**

Convert an object of class 'model' into a ready-to-write vector of string

**Usage**

```
## S3 method for class 'model'  
toString(x, ...)
```

**Arguments**

- x model object, object to be converted.
- ... Optional arguments.

**Value**

A string ready to be printed or written.

**Examples**

```
toString(model())
```

---

toString.parameter     *parameter to string*

---

### Description

Convert an object of class 'parameter' into a ready-to-write vector of string

### Usage

```
## S3 method for class 'parameter'  
toString(x, ...)
```

### Arguments

x                parameter object, object to be converted.  
...                Optional arguments.

### Value

A string ready to be printed or written.

### Examples

```
p <- parameter(name='par',init=0,prior.dist='Gaussian',prior.par=c(0,1))  
toString(p)
```

---

toString.parameter\_VAR  
                        *parameter\_VAR to string*

---

### Description

Convert an object of class 'parameter\_VAR' into a ready-to-write vector of string

### Usage

```
## S3 method for class 'parameter_VAR'  
toString(x, ...)
```

### Arguments

x                parameter\_VAR object, object to be converted.  
...                Optional arguments.

**Value**

A string ready to be printed or written.

**Examples**

```
X=data.frame(input1=rnorm(100),input2=rnorm(100))
Y=data.frame(output=X$input1+0.8*X$input2+0.1*rnorm(100))
VAR.indx=data.frame(indx=c(rep(1,50),rep(2,50)))
workspace=tempdir()
d <- dataset(X=X,Y=Y,data.dir=workspace,VAR.indx=VAR.indx)
p <- parameter_VAR(name='par',index='indx',d=d,
                      init=c(-1,1,2),
                      prior.dist=c('Gaussian','FlatPrior','Triangle'),
                      prior.par=list(c(-1,1),NULL,c(2,0,5)))
toString(p)
```

**toString.prediction** *prediction to string*

**Description**

Convert an object of class 'prediction' into a ready-to-write vector of string

**Usage**

```
## S3 method for class 'prediction'
toString(x, ...)
```

**Arguments**

x	prediction object, object to be converted.
...	Optional arguments.

**Value**

A string ready to be printed or written.

**Examples**

```
pred=prediction(X=twoPopulations[,1:3],spagFiles=c('P1.spag','P2.spag'))
toString(pred)
```

---

toString.remnantErrorModel  
*remnantErrorModel to string*

---

## Description

Convert an object of class 'remnantErrorModel' into a ready-to-write vector of string

## Usage

```
## S3 method for class 'remnantErrorModel'  
toString(x, ...)
```

## Arguments

x               remnantErrorModel object, object to be converted.  
...              Optional arguments.

## Value

A string ready to be printed or written.

## Examples

```
toString(remnantErrorModel())
```

---

toString.residualOptions  
*residualOptions to string*

---

## Description

Convert an object of class 'residualOptions' into a ready-to-write vector of string

## Usage

```
## S3 method for class 'residualOptions'  
toString(x, ...)
```

## Arguments

x               residualOptions object, object to be converted.  
...              Optional arguments.

**Value**

A string ready to be printed or written.

**Examples**

```
toString(residualOptions())
```

<code>toString.runOptions</code>	<i>runOptions to string</i>
----------------------------------	-----------------------------

**Description**

Convert an object of class 'runOptions' into a ready-to-write vector of string

**Usage**

```
## S3 method for class 'runOptions'  
toString(x, ...)
```

**Arguments**

<code>x</code>	runOptions object, object to be converted.
<code>...</code>	Optional arguments.

**Value**

A string ready to be printed or written.

**Examples**

```
toString(runOptions())
```

<code>tracePlot</code>	<i>MCMC reporting</i>
------------------------	-----------------------

**Description**

2DO (adapt from STooDs): Generate pdf report files summarizing mcmc samples

**Usage**

```
tracePlot(sim, ylab = "values", keep = NULL, col = "black", psize = 0.5)
```

### Arguments

sim	vector or matrix or data frame, MCMC simulations
ylab	Character, label of y-axis to be used if sim has no names
keep	Integer vector, indices of samples to be kept in cooked MCMC sample
col	Color
psize	Numeric, point size

### Details

tracePlot

returns a trace plot ggplot (or a list thereof if several columns in sim)

### Value

A ggplot (or a list thereof if several columns in sim)

### Examples

```
# Create Monte Carlo samples
n=1000
sim=data.frame(p1=rnorm(n),p2=rlnorm(n),p3=runif(n))
# create trace plot for each component
figures=tracePlot(sim)
```

twoPopulations      *Evolution of two populations*

### Description

Size of two populations of the same species put in two different environments, as a function of time and local temperature.

- Input variables: time t, temperature at site 1 T1, temperature at site 2 T2.
- Output variables: population size at site 1 P1, population size at site 2 P2.
- Data are synthetically generated from a logistic model.

### Usage

twoPopulations

### Format

An object of class `data.frame` with 101 rows and 5 columns.

**violinPlot***violinPlot***Description**

returns a violinplot ggplot

**Usage**

```
violinPlot(sim, ylab = "values", col = "black")
```

**Arguments**

sim	vector or matrix or data frame, MCMC simulations
ylab	Character, label of y-axis
col	Color

**Value**

A ggplot

**Examples**

```
# Create Monte Carlo samples
n=1000
sim=data.frame(p1=rnorm(n),p2=rlnorm(n),p3=runif(n))
# create violin plot comparing all components
figure=violinPlot(sim)
```

**writePredInputs***Write prediction inputs***Description**

Write input data of the prediction into files

**Usage**

```
writePredInputs(o)
```

**Arguments**

o	prediction object
---	-------------------

**Value**

nothing - just write to files.

## Examples

```
temp=tempdir()
pred=prediction(X=twoPopulations[,1:3],spagFiles=c('P1.spag','P2.spag'),
                 data.dir=temp)
writePredInputs(pred)
```

---

xtraModelInfo      *xtraModelInfo constructor.*

---

## Description

Creates a new instance of a 'xtraModelInfo' object containing extra model information

## Usage

```
xtraModelInfo(fname = "Config_Xtra.txt", object = NULL)
```

## Arguments

fname	Character, configuration file name.
object	any R object containing xtra model info - typically a list of stuff. The content and meaning of 'object' is completely model-specific.

## Value

An object of class 'xtraModelInfo'.

## Examples

```
x <- xtraModelInfo()
```

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