

# Package ‘JointAI’

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**Version** 0.5.0

**Title** Joint Analysis and Imputation of Incomplete Data

**Description** Provides joint analysis and imputation of (generalized) linear and cumulative logit regression models, (generalized) linear and cumulative logit mixed models and parametric (Weibull) as well as Cox proportional hazards survival models with incomplete (covariate) data in the Bayesian framework.  
The package performs some preprocessing of the data and creates a 'JAGS' model, which will then automatically be passed to 'JAGS' <<http://mcmc-jags.sourceforge.net>> with the help of the package 'rjags'.  
It also provides summary and plotting functions for the output and allows to export imputed values.

**URL** <https://nerler.github.io/JointAI>

**License** GPL (>= 2)

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**BugReports** <https://github.com/nerler/JointAI/issues>

**LazyData** TRUE

**RoxygenNote** 6.1.1

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add_samples	<i>Add samples to an object of class JointAI</i>
-------------	--

---

### Description

Allows to continue sampling from an existing object of class 'JointAI'. When the original sample was created using parallel computation, the separate 'jags' objects will be recompiled and sampling will again be performed in parallel.

### Usage

```
add_samples(object, n.iter, add = TRUE, thin = NULL,
            monitor_params = NULL, progress.bar = "text", mess = TRUE)
```

**Arguments**

object	object inheriting from class 'JointAI'
n.iter	number of iterations to monitor
add	logical; should the new MCMC samples be added to the existing samples or replace them? If samples are added, thin and var.names are ignored.
thin	thinning interval (see <a href="#">window.mcmc</a> )
monitor_params	named vector specifying which parameters should be monitored
progress.bar	character string specifying the type of progress bar. Possible values are "text", "gui", and "none". See <a href="#">update</a> .
mess	logical; should messages be given? Default is TRUE. Note: this applies only to messages given directly by <b>JointAI</b> .

**See Also**

[lm\\_imp](#), [glm\\_imp](#), [lme\\_imp](#), [clm\\_imp](#), [glme\\_imp](#), [clmm\\_imp](#), [survreg\\_imp](#), [coxph\\_imp](#)

The vignette [Parameter Selection](#) contains some examples how to specify the argument monitor\_params.

**Examples**

```
# Example 1:
# Run an initial JointAI model:
mod <- lm_imp(y~C1 + C2 + M2, data = wideDF, n.iter = 100)

# Continue sampling:
mod_add <- add_samples(mod, n.iter = 200, add = TRUE)

# Example 2:
# Continue sampling, but additionally sample imputed values.
# Note: Setting different parameters to monitor than in the original model
# requires add = FALSE.
imps <- add_samples(mod, n.iter = 200, monitor_params = c("imps" = TRUE),
                    add = FALSE)
```

---

default\_hyperpars      *Get default values for hyperparameters*

---

**Description**

Prints the list of default values for the hyperparameters.

**Usage**

```
default_hyperpars()
```

**Details****norm:** hyperparameters for normal and lognormal models

mu_reg_norm	mean in the priors for regression coefficients
tau_reg_norm	precision in the priors for regression coefficients
shape_tau_norm	shape parameter in Gamma prior for precision of imputed variable
rate_tau_norm	rate parameter in Gamma prior for precision of imputed variable

**gamma:** hyperparameters for Gamma models

mu_reg_gamma	mean in the priors for regression coefficients
tau_reg_gamma	precision in the priors for regression coefficients
shape_tau_gamma	shape parameter in Gamma prior for precision of imputed variable
rate_tau_gamma	rate parameter in Gamma prior for precision of imputed variable

**beta:** hyperparameters for beta models

mu_reg_beta	mean in the priors for regression coefficients
tau_reg_beta	precision in the priors for regression coefficients
shape_tau_beta	shape parameter in Gamma prior for precision of imputed variable
rate_tau_beta	rate parameter in Gamma prior for precision of imputed variable

**logit:** hyperparameters for logistic models

mu_reg_logit	mean in the priors for regression coefficients
tau_reg_logit	precision in the priors for regression coefficients

**probit:** hyperparameters for probit models

mu_reg_logit	mean in the priors for regression coefficients
tau_reg_logit	precision in the priors for regression coefficients

**multinomial:** hyperparameters for multinomial models

mu_reg_multinomial	mean in the priors for regression coefficients
tau_reg_multinomial	precision in the priors for regression coefficients

**ordinal:** hyperparameters for ordinal models

mu_reg_ordinal	mean in the priors for regression coefficients
tau_reg_ordinal	precision in the priors for regression coefficients
mu_delta_ordinal	mean in the prior for the intercepts
tau_delta_ordinal	precision in the priors for the intercepts

**Z:** function creating hyperparameters for the random effects in mixed models, with output elements

RinvD	scale matrix in Wishart prior (*) for random effects covariance matrix
KinvD	degrees of freedom in Wishart prior for random effects covariance matrix
shape_diag_RinvD	shape parameter in Gamma prior for the diagonal elements of RinvD
rate_diag_RinvD	rate parameter in Gamma prior for the diagonal elements of RinvD

(\*) when there is only one random effect a Gamma distribution is used instead of the Wishart and RinvD and KinvD are NULL

**surv:** parameters for survival models (parametric and proportional hazard)

mu_reg_surv	mean in the priors for regression coefficients
tau_reg_surv	precision in the priors for regression coefficients

**coxph:** parameters for Cox proportional hazards models

c	confidence in prior guess for the hazard function
r	failure rate per unit time
eps	time increment

## Examples

```
default_hyperpars()
```

---

densplot	<i>Plot posterior density from JointAI model</i>
----------	--

---

## Description

Plots a set of densities (per MC chain and coefficient) from the MCMC sample of an object of class "JointAI".

## Usage

```
densplot(object, ...)

## S3 method for class 'mcmc.list'
densplot(object, start = NULL, end = NULL,
  thin = NULL, ...)

## S3 method for class 'JointAI'
densplot(object, start = NULL, end = NULL,
  thin = NULL, subset = c(analysis_main = TRUE), vlines = NULL,
  nrow = NULL, ncol = NULL, joined = FALSE, use_ggplot = FALSE,
  keep_aux = FALSE, warn = TRUE, ...)
```

**Arguments**

object	object inheriting from class 'JointAI'
...	additional parameters passed to <code>plot</code>
start	the first iteration of interest (see <code>window.mcmc</code> )
end	the last iteration of interest (see <code>window.mcmc</code> )
thin	thinning interval (see <code>window.mcmc</code> )
subset	subset of parameters/variables/nodes (columns in the MCMC sample). Uses the same logic as the argument <code>monitor_params</code> in <code>lm_imp</code> , <code>glm_imp</code> , <code>clm_imp</code> , <code>lme_imp</code> , <code>glme_imp</code> , <code>survreg_imp</code> and <code>coxph_imp</code> .
vlines	list, where each element is a named list of parameters that can be passed to <code>abline</code> to create vertical lines. Each of the list elements needs to contain at least <code>v = &lt;x location&gt;</code> , where <code>&lt;x location&gt;</code> is a vector of the same length as the number of plots (see examples).
nrow	optional number of rows and columns in the plot layout; automatically chosen if unspecified
ncol	optional number of rows and columns in the plot layout; automatically chosen if unspecified
joined	logical; should the chains be combined before plotting?
use_ggplot	logical; Should ggplot be used instead of the base graphics?
keep_aux	logical; Should constant effects of auxiliary variables be kept in the output?
warn	logical; should warnings be given? Default is TRUE. Note: this applies only to warnings given directly by <b>JointAI</b> .

**See Also**

The vignette [Parameter Selection](#) contains some examples how to specify the argument `subset`.

**Examples**

```
# fit a JointAI object:
mod <- lm_imp(y ~ C1 + C2 + M2, data = wideDF, n.iter = 100)

# Example 1: basic densityplot
densplot(mod)

# Example 2: use vlines to mark zero
densplot(mod, col = c("darkred", "darkblue", "darkgreen"),
          vlines = list(list(v = rep(0, nrow(summary(mod)$stats)),
                           col = grey(0.8))))

# Example 3: use vlines to visualize the posterior mean and 2.5% and 97.5% quantiles
densplot(mod, vlines = list(list(v = summary(mod)$stats[, "Mean"], lty = 1, lwd = 2),
                            list(v = summary(mod)$stats[, "2.5%"], lty = 2),
```

```

list(v = summary(mod)$stats[, "97.5%"], lty = 2)))

# Example 4: ggplot version
densplot(mod, use_ggplot = TRUE)

# Example 5: changing how the ggplot version looks (using standard ggplot syntax)
library(ggplot2)

densplot(mod, use_ggplot = TRUE) +
  xlab("value") +
  theme(legend.position = 'bottom') +
  scale_color_brewer(palette = 'Dark2', name = 'chain')

```

---

get\_MIdat

*Extract multiple imputed datasets*


---

### Description

Creates a dataset containing multiple imputed datasets stacked onto each other (i.e., long format). These data can be automatically exported to SPSS (i.e., a .txt file containing the data and a .sps file containing syntax to generate a .sav file). For the export function the **foreign** package needs to be installed.

### Usage

```

get_MIdat(object, m = 10, include = TRUE, start = NULL,
  minspace = 50, seed = NULL, export_to_SPSS = FALSE,
  resdir = NULL, filename = NULL)

```

### Arguments

object	object inheriting from class 'JointAI'
m	number of imputed datasets
include	should the original, incomplete data be included?
start	the first iteration of interest (see <a href="#">window.mcmc</a> )
minspace	minimum number of iterations between iterations chosen as imputed values.
seed	optional seed
export_to_SPSS	logical; should the completed data be exported to SPSS?
resdir	optional directory for results (if unspecified and export_to_SPSS = TRUE the current working directory is used)
filename	optional file name (without ending; if unspecified and export_to_SPSS = TRUE a name is generated automatically)

**Value**

A dataframe containing the imputed values (and original data) stacked. The variable `Imputation_` identifies the imputations, while `.rownr` identifies rows of the rows of the original data. In cross-sectional datasets the variable `.id` is added as subject identifier.

**See Also**

[plot\\_imp\\_distr](#)

**Examples**

```
# fit a model and monitor the imputed values with monitor_params = c(imps = TRUE)
mod <- lm_imp(y~C1 + C2 + M2, data = wideDF, monitor_params = c(imps = TRUE), n.iter = 100)

# Example 1: without export to SPSS
MIs <- get_MIdat(mod, m = 3, seed = 123)

## Not run:
# Example 2: with export for SPSS (here: to the temporary directory "temp_dir")
temp_dir <- tempdir()
MIs <- get_MIdat(mod, m = 3, seed = 123, resdir = temp_dir,
                filename = "example_imputation",
                export_to_SPSS = TRUE)

## End(Not run)
```

---

get\_models

*Set default (imputation) models and order*

---

**Description**

Set default (imputation) models and order

**Usage**

```
get_models(fixed, random = NULL, data, auxvars = NULL,
           no_model = NULL)

get_imp_meth(fixed, random = NULL, data, auxvars = NULL,
             no_model = NULL)
```

**Arguments**

fixed	a two sided formula describing the fixed-effects part of the model (see <a href="#">formula</a> )
random	only for <code>lme_imp</code> , <code>glme_imp</code> and <code>clmm_imp</code> : a one-sided formula of the form $\sim x_1 + \dots + x_n \mid g$ , where $x_1 + \dots + x_n$ specifies the model for the random effects and <code>g</code> the grouping variable



data	a data.frame
auxvars	optional vector of variable names that should be used as predictors in the imputation procedure (and will be imputed if necessary) but are not part of the analysis model
no_model	names of variables for which no model should be specified. Note that this is only possible for completely observed variables and may imply assumptions of independence between the excluded variable and incomplete variables.

### Value

`get_models()` returns a list of two vectors named `models` and `meth`. `models` is a named vector containing the names of covariates that either have missing values and/or are longitudinal (level-1) covariates and the corresponding default (imputation) models. `meth` is a subset of `models` containing only the variables that have missing values.

### Examples

```
get_models(y ~ C1 + C2 + B2 + O2 + M2, data = wideDF)
get_imp_meth(y ~ C1 + C2 + B2 + O2 + M2, data = wideDF)

get_models(y ~ C1 + O2 + c2 + b1 + o2 + time, random = ~ 1 | id, data = longDF)
get_imp_meth(y ~ C1 + O2 + c2 + b1 + o2 + time, random = ~ 1 | id, data = longDF)

get_models(y ~ C1 + O2 + c2 + b1 + o2 + time, random = ~ 1 | id,
           no_model = 'time', data = longDF)
```

---

GR\_crit

*Gelman-Rubin criterion for convergence*


---

### Description

Calculates the Gelman-Rubin criterion for convergence (uses [gelman.diag](#) from package `coda`).

### Usage

```
GR_crit(object, confidence = 0.95, transform = FALSE,
        autoburnin = TRUE, multivariate = TRUE, subset = NULL,
        start = NULL, end = NULL, thin = NULL, warn = TRUE, ...)
```

### Arguments

object	object inheriting from class 'JointAI'
confidence	the coverage probability of the confidence interval for the potential scale reduction factor
transform	a logical flag indicating whether variables in <code>x</code> should be transformed to improve the normality of the distribution. If set to <code>TRUE</code> , a log transform or logit transform, as appropriate, will be applied.

autoburnin	a logical flag indicating whether only the second half of the series should be used in the computation. If set to TRUE (default) and <code>start(x)</code> is less than <code>end(x)/2</code> then start of series will be adjusted so that only second half of series is used.
multivariate	a logical flag indicating whether the multivariate potential scale reduction factor should be calculated for multivariate chains
subset	subset of parameters/variables/nodes (columns in the MCMC sample). Uses the same logic as the argument <code>monitor_params</code> in <code>lm_imp</code> , <code>glm_imp</code> , <code>clm_imp</code> , <code>lme_imp</code> , <code>glme_imp</code> , <code>survreg_imp</code> and <code>coxph_imp</code> .
start	the first iteration of interest (see <code>window.mcmc</code> )
end	the last iteration of interest (see <code>window.mcmc</code> )
thin	thinning interval (see <code>window.mcmc</code> )
warn	logical; should warnings be given? Default is TRUE. Note: this applies only to warnings given directly by <b>JointAI</b> .
...	currently not used

## References

Gelman, A and Rubin, DB (1992) Inference from iterative simulation using multiple sequences, *Statistical Science*, **7**, 457-511.

Brooks, SP. and Gelman, A. (1998) General methods for monitoring convergence of iterative simulations. *Journal of Computational and Graphical Statistics*, **7**, 434-455.

## See Also

The vignette [Parameter Selection](#) contains some examples how to specify the argument `subset`.

## Examples

```
mod1 <- lm_imp(y~C1 + C2 + M2, data = wideDF, n.iter = 100)
GR_crit(mod1)
```

## Description

The **JointAI** package performs simultaneous imputation and inference for incomplete data using the Bayesian framework. Distributions of incomplete variables, conditional on other covariates, are specified automatically and modeled jointly with the analysis model. MCMC sampling is performed in '**JAGS**' via the R package **rjags**.

## Main functions

The package has the following main functions that allow analysis in different settings:

- `lm_imp` for linear regression
- `glm_imp` for generalized linear regression
- `clm_imp` for cumulative logit models
- `lme_imp` for linear mixed models
- `glme_imp` for generalized linear mixed models
- `clmm_imp` for cumulative logit mixed models
- `survreg_imp` for parametric (Weibull) survival models
- `coxph_imp` for Cox proportional hazard models

As far as possible, the specification of these functions is analogue to the specification of their complete data versions `lm`, `glm`, `clm` (from the package **ordinal**), `lme` (from the package **nlme**), `clmm2` (from the package **ordinal**), `survreg` (from the package **survival**) and `coxph` (from the package **survival**).

Results can be summarized and printed with `summary.JointAI`, `coef.JointAI` and `confint.JointAI`, and visualized using `traceplot` or `densplot`. The function `predict.JointAI` allows prediction (including credible intervals) from JointAI models.

## Evaluation and export

Two criteria for evaluation of convergence and precision of the posterior estimate are available:

- `GR_crit` implements the Gelman-Rubin criterion ('potential scale reduction factor') for convergence
- `MC_error` calculates the Monte Carlo error to evaluate the precision of the MCMC sample

Imputed data can be extracted (and exported to SPSS) using `get_MIdat`. The function `plot_imp_distr` allows visual comparison of the distribution of observed and imputed values.

## Other useful functions

- `parameters` and `list_impmodels` to gain insight in the specified model
- `plot_all` and `md_pattern` to visualize the distribution of the data and the missing data pattern

## Vignettes

The following vignettes are available

- *Minimal Example:*  
A minimal example demonstrating the use of `lm_imp`, `summary.JointAI`, `traceplot` and `densplot`.
- *Visualizing Incomplete Data:*  
Demonstrations of the options in `plot_all` (plotting histograms and barplots for all variables in the data) and `md_pattern` (plotting or printing the missing data pattern).

- **Model Specification:**  
Explanation and demonstration of all parameters that are required or optional to specify the model structure in `lm_imp`, `glm_imp` and `lme_imp`. Among others, the functions `parameters`, `list_impmodels`, `get_imp_meth` and `set_refcat` are used.
- **Parameter Selection:**  
Examples on how to select the parameters/variables/nodes to follow using the argument `monitor_params` and the parameters/variables/nodes displayed in the `summary`, `traceplot`, `densplot` or when using `GR_crit` or `MC_error`.
- **MCMC Settings:**  
Examples demonstrating how to set the arguments controlling settings of the MCMC sampling, i.e., `n.adapt`, `n.iter`, `n.chains`, `thin`, `inits`.
- **After Fitting:**  
Examples on the use of functions to be applied after the model has been fitted, including `traceplot`, `densplot`, `summary`, `GR_crit`, `MC_error`, `predict`, `predDF` and `get_MIdat`.

## References

Erler, N.S., Rizopoulos, D., Rosmalen, J., Jaddoe, V.W.V., Franco, O. H., & Lesaffre, E.M.E.H. (2016). Dealing with missing covariates in epidemiologic studies: A comparison between multiple imputation and a full Bayesian approach. *Statistics in Medicine*, 35(17), 2955-2974. doi: [10.1002/sim.6944](https://doi.org/10.1002/sim.6944)

Erler, N.S., Rizopoulos D., Jaddoe, V.W.V., Franco, O.H. & Lesaffre, E.M.E.H. (2019). Bayesian imputation of time-varying covariates in linear mixed models. *Statistical Methods in Medical Research*, 28(2), 555–568. doi: [10.1177/0962280217730851](https://doi.org/10.1177/0962280217730851)

---

JointAIObject

*Fitted object of class JointAI*

---

## Description

An object returned by one of the functions `lm_imp()`, `glm_imp()`, `clm_imp()`, `lme_imp()`, `glme_imp()`, `clmm_imp()`, `survreg_imp()` or `coxph_imp()`.

## Value

<code>analysis_type</code>	<code>lm</code> , <code>glm</code> , <code>clm</code> , <code>lme</code> , <code>glme</code> , <code>clmm</code> , <code>survreg</code> or <code>coxph</code> with attributes <code>family</code> and <code>link</code>
<code>data</code>	the original (incomplete) dataset
<code>models</code>	named vector specifying the models used for longitudinal and incomplete covariates
<code>fixed</code>	supplied fixed effects formula
<code>random</code>	supplied random effects formula
<code>Mlist</code>	a list: containing the data, split up into <ul style="list-style-type: none"> <li>• <code>outcome (y)</code></li> </ul>

- censoring indicator for survival outcomes (*cens*)
- cross-sectional main effects (*Xc*)
- cross-sectional interactions (*Xic*)
- longitudinal main effects (*Xl*)
- longitudinal interactions (*Xil*)
- categorical cross-sectional incomplete variables (*Xcat*)
- categorical longitudinal variables (*Xlcat*)
- transformed cross-sectional variables (*Xtrafo*)
- random effects design matrix (*Z*)

and other important specifications:

- a list naming which columns of the above matrices are covariates in the analysis model (*cols\_main*)
- a list giving the names of the covariates in the analysis model per matrix (*names\_main*)
- specification for transformations (*trafos*)
- specification for hierarchical centering (*hc\_list*)
- reference values and dummies for categorical variables (*refs*)
- vector of auxiliary variables (*auxvars*)
- grouping specification (*groups*)
- the vector of variables to be scaled (*scale\_vars*)
- updated fixed effects structure (*fixed2*)
- the number of categories if the outcome of the analysis model is categorical (*ncat*)
- the number of subjects (*N*)
- whether posterior predictive checks are to be enabled *ppc* (not yet used)
- whether ridge shrinkage priors should be used for the regression coefficients of the analysis model (*ridge*)
- the number of random effects (*nraneff*)

<i>K</i>	matrix specifying the indices of the regression coefficients that are related to different parts of the model
<i>K_imp</i>	matrix specifying the indices of regression coefficients for the imputation models relating to different covariates
<i>mcmc_settings</i>	a list with elements <ul style="list-style-type: none"> <li><i>modelfile</i> name and path of JAGS model file</li> <li><i>n.chains</i> number of MCMC chains</li> <li><i>n.adapt</i> number of iterations in the adaptive phase</li> <li><i>n.iter</i> number of iterations in the MCMC sample</li> <li><i>variable.names</i> monitored nodes</li> <li><i>thin</i> thinning of the MCMC sample</li> <li><i>inits</i> a list containing the initial values that were passed to <b>rjags</b></li> <li><i>parallel</i> whether parallel sampling was used</li> <li><i>ncores</i> how many cores were used in parallel sampling</li> </ul>

monitor_params	the list of parameter groups to be monitored
data_list	list with data that was passed to <b>rjags</b>
scale_pars	matrix with parameters used to center and scale the continuous variables
hyperpars	a list containing the values of the hyperparameters used
imp_par_list	a list with parameters used to write the imputation model syntax
model	JAGS model
sample	MCMC sample on the sampling scale (included only if keep_scaled_sample = TRUE)
MCMC	MCMC sample, scaled back to the scale of the data
time	the computational time used for the sampling (adaptive phase + sampling)
call	the original call

---

list_models	<i>List imputation models</i>
-------------	-------------------------------

---

### Description

Print information on all models for incomplete covariates used in a JointAI object, including the model type, names of the parameters used and hyperparameters.

### Usage

```
list_models(object, predvars = TRUE, regcoef = TRUE,
            otherpars = TRUE, priors = TRUE, refcat = TRUE)
```

```
list_impmodels(object, predvars = TRUE, regcoef = TRUE,
               otherpars = TRUE, priors = TRUE, refcat = TRUE)
```

### Arguments

object	object inheriting from class 'JointAI'
predvars	logical; should information on the predictor variables be printed?
regcoef	logical; should information on the regression coefficients be printed?
otherpars	logical; should information on other parameters be printed?
priors	logical; should information on the priors be printed?
refcat	logical; should information on the reference category be printed?

**Note**

The models listed by this function are not the actual imputation models, but the conditional models that are part of the specification of the joint distribution of the data. Briefly, the joint distribution is specified as a sequence of conditional models

$$p(y|x_1, x_2, x_3, \dots, \theta)p(x_1|x_2, x_3, \dots, \theta)p(x_2|x_3, \dots, \theta)\dots$$

The actual imputation models are the full conditional distributions  $p(x_1|\cdot)$  derived from this joint distribution. Even though the conditional distributions do not contain the outcome and all other covariates in their linear predictor, outcome and other covariates are taken into account implicitly, since imputations are sampled from the full conditional distributions. For more details, see Erler et al. (2016).

The function `list_models` prints information on the conditional distributions of the incomplete covariates (since they are what is specified; the full-conditionals are automatically derived within JAGS). The outcome is, thus, not part of the printed linear predictor, but is still included during imputation.

**References**

Erler, N. S., Rizopoulos, D., Rosmalen, J. V., Jaddoe, V. W., Franco, O. H., & Lesaffre, E. M. (2016). Dealing with missing covariates in epidemiologic studies: A comparison between multiple imputation and a full Bayesian approach. *Statistics in Medicine*, 35(17), 2955-2974.

**Examples**

```
# (set n.adapt = 0 and n.iter = 0 to prevent MCMC sampling to save time)
mod1 <- lm_imp(y ~ C1 + C2 + M2 + O2 + B2, data = wideDF, n.adapt = 0, n.iter = 0)

list_models(mod1)
```

---

longDF

*Longitudinal example dataset*


---

**Description**

A simulated longitudinal dataset.

**Usage**

```
data(longDF)
```

**Format**

A simulated data frame with 329 rows and 21 variables with data from 100 subjects:

**C1** continuous, complete baseline variable  
**C2** continuous, incomplete baseline variable  
**B1** binary, complete baseline variable  
**B2** binary, incomplete baseline variable  
**M1** unordered factor; complete baseline variable  
**M2** unordered factor; incomplete baseline variable  
**O1** ordered factor; complete baseline variable  
**O2** ordered factor; incomplete baseline variable  
**P1** count variable; complete baseline variable  
**P2** count variable; incomplete baseline variable  
**c1** continuous, complete longitudinal variable  
**c2** continuous incomplete longitudinal variable  
**b1** binary, complete longitudinal variable  
**b2** binary incomplete longitudinal variable  
**o1** ordered factor; complete longitudinal variable  
**o2** ordered factor; incomplete longitudinal variable  
**p1** count variable; complete longitudinal variable  
**p2** count variable; incomplete longitudinal variable  
**id** id (grouping) variable  
**time** continuous complete longitudinal variable  
**y** continuous, longitudinal (outcome) variable

---

MC\_error

*Monte Carlo error*

---

**Description**

Calculate and plot the Monte Carlo error of the samples from a JointAI model.

**Usage**

```
MC_error(x, subset = NULL, start = NULL, end = NULL, thin = NULL,
         digits = 2, warn = TRUE, ...)

## S3 method for class 'MCElist'
plot(x, data_scale = TRUE, plotpars = NULL,
     ablinepars = list(v = 0.05), ...)
```



**Arguments**

x	object inheriting from class 'JointAI'
subset	subset of parameters/variables/nodes (columns in the MCMC sample). Uses the same logic as the argument <code>monitor_params</code> in <code>lm_imp</code> , <code>glm_imp</code> , <code>clm_imp</code> , <code>lme_imp</code> , <code>glme_imp</code> , <code>survreg_imp</code> and <code>coxph_imp</code> .
start	the first iteration of interest (see <code>window.mcmc</code> )
end	the last iteration of interest (see <code>window.mcmc</code> )
thin	thinning interval (see <code>window.mcmc</code> )
digits	number of digits for output
warn	logical; should warnings be given? Default is TRUE. Note: this applies only to warnings given directly by <b>JointAI</b> .
...	Arguments passed on to <code>mcmcse:mcse.mat</code>
	<b>size</b> the batch size. The default value is "sqrt", which uses the square root of the sample size. "cuberoot" will cause the function to use the cube root of the sample size. A numeric value may be provided if neither "sqrt" nor "cuberoot" is satisfactory.
	<b>g</b> a function such that $E(g(x))$ is the quantity of interest. The default is NULL, which causes the identity function to be used.
	<b>method</b> the method used to compute the standard error. This is one of "bm" (batch means, the default), "obm" (overlapping batch means), "tukey" (spectral variance method with a Tukey-Hanning window), or "bartlett" (spectral variance method with a Bartlett window).
data_scale	show the Monte Carlo error of the sample transformed back to the scale of the data (TRUE) or on the sampling scale (this requires the argument <code>keep_scaled_mcmc = TRUE</code> in the JointAI model)
plotpars	optional; list of parameters passed to <code>plot()</code>
ablinepars	optional; list of parameters passed to <code>abline()</code>

**Value**

An object of class `MCElist` with elements `unscaled`, `scaled` and `digits`. The first two are matrices with columns `est` (posterior mean), `MCSE` (Monte Carlo error), `SD` (posterior standard deviation) and `MCSE/SD` (Monte Carlo error divided by post. standard deviation.)

**Methods (by generic)**

- `plot`: plot Monte Carlo error

**Note**

Lesaffre & Lawson (2012) [p. 195] suggest the Monte Carlo error of a parameter should not be more than 5% of the posterior standard deviation of this parameter (i.e.,  $MCSE/SD \leq 0.05$ ).

**References**

Lesaffre, E., & Lawson, A. B. (2012). *Bayesian Biostatistics*. John Wiley & Sons.

**See Also**

The vignette [Parameter Selection](#) contains some examples how to specify the argument subset.

**Examples**

```
mod <- lm_imp(y~C1 + C2 + M2, data = wideDF, n.iter = 100)

MC_error(mod)

plot(MC_error(mod), ablinepars = list(lty = 2))
```

---

 md\_pattern

*Missing data pattern*


---

**Description**

Plot the pattern of missing data.

**Usage**

```
md_pattern(data, color = c(grDevices::grey(0.1), grDevices::grey(0.7)),
  border = grDevices::grey(0.5), plot = TRUE, pattern = FALSE,
  print_xaxis = TRUE, ylab = "Number of observations per pattern",
  print_yaxis = TRUE, legend.position = "bottom", ...)
```

**Arguments**

data	data frame
color	vector of length two, that specifies the color used to indicate observed and missing (in that order)
border	color of the grid
plot	logical; should the missing data pattern be plotted?
pattern	logical; should the missing data pattern be returned as matrix?
print_xaxis, print_yaxis	logical; should the x-axis (below the plot) and y-axis (on the right) be printed?
ylab	y-axis label
legend.position	the position of legends ("none", "left", "right", "bottom", "top", or two-element numeric vector)
...	optional additional parameters, currently not used

**Note**

This function requires the [ggplot2](#) package to be installed.

**See Also**

See the vignette [Visualizing Incomplete Data](#) for more examples.

**Examples**

```
par(mar = c(3, 1, 1.5, 1.5), mgp = c(2, 0.6, 0))
md_pattern(wideDF)
```

---

 model\_imp

*Joint analysis and imputation of incomplete data*


---

**Description**

Functions to estimate (generalized) linear and (generalized) linear mixed models, ordinal and ordinal mixed models, and parametric (Weibull) as well as Cox proportional hazards survival models using MCMC sampling, while imputing missing values.

**Usage**

```
lm_imp(formula, data, n.chains = 3, n.adapt = 100, n.iter = 0,
  thin = 1, monitor_params = NULL, auxvars = NULL, refcats = NULL,
  models = NULL, no_model = NULL, trunc = NULL, ridge = FALSE,
  ppc = TRUE, seed = NULL, inits = NULL, parallel = FALSE,
  ncores = NULL, scale_vars = NULL, scale_pars = NULL,
  hyperpars = NULL, modelname = NULL, modeldir = NULL,
  keep_model = FALSE, overwrite = NULL, quiet = TRUE,
  progress.bar = "text", warn = TRUE, mess = TRUE,
  keep_scaled_mcmc = FALSE, ...)
```

```
glm_imp(formula, family, data, n.chains = 3, n.adapt = 100,
  n.iter = 0, thin = 1, monitor_params = NULL, auxvars = NULL,
  refcats = NULL, models = NULL, no_model = NULL, trunc = NULL,
  ridge = FALSE, ppc = TRUE, seed = NULL, inits = NULL,
  parallel = FALSE, ncores = NULL, scale_vars = NULL,
  scale_pars = NULL, hyperpars = NULL, modelname = NULL,
  modeldir = NULL, keep_model = FALSE, overwrite = NULL,
  quiet = TRUE, progress.bar = "text", warn = TRUE, mess = TRUE,
  keep_scaled_mcmc = FALSE, ...)
```

```
clm_imp(fixed, data, n.chains = 3, n.adapt = 100, n.iter = 0,
  thin = 1, monitor_params = NULL, auxvars = NULL, refcats = NULL,
  models = NULL, no_model = NULL, trunc = NULL, ridge = FALSE,
  ppc = TRUE, seed = NULL, inits = NULL, parallel = FALSE,
  ncores = NULL, scale_vars = NULL, scale_pars = NULL,
  hyperpars = NULL, modelname = NULL, modeldir = NULL,
```

```
keep_model = FALSE, overwrite = NULL, quiet = TRUE,  
progress.bar = "text", warn = TRUE, mess = TRUE,  
keep_scaled_mcmc = FALSE, ...)
```

```
lme_imp(fixed, data, random, n.chains = 3, n.adapt = 100, n.iter = 0,  
  thin = 1, monitor_params = NULL, auxvars = NULL, refcats = NULL,  
  models = NULL, no_model = NULL, trunc = NULL, ridge = FALSE,  
  ppc = TRUE, seed = NULL, inits = NULL, parallel = FALSE,  
  ncores = NULL, scale_vars = NULL, scale_pars = NULL,  
  hyperpars = NULL, modelname = NULL, modeldir = NULL,  
  keep_model = FALSE, overwrite = NULL, quiet = TRUE,  
  progress.bar = "text", warn = TRUE, mess = TRUE,  
  keep_scaled_mcmc = FALSE, ...)
```

```
glme_imp(fixed, data, random, family, n.chains = 3, n.adapt = 100,  
  n.iter = 0, thin = 1, monitor_params = NULL, auxvars = NULL,  
  refcats = NULL, models = NULL, no_model = NULL, trunc = NULL,  
  ridge = FALSE, ppc = TRUE, seed = NULL, inits = NULL,  
  parallel = FALSE, ncores = NULL, scale_vars = NULL,  
  scale_pars = NULL, hyperpars = NULL, modelname = NULL,  
  modeldir = NULL, keep_model = FALSE, overwrite = NULL,  
  quiet = TRUE, progress.bar = "text", warn = TRUE, mess = TRUE,  
  keep_scaled_mcmc = FALSE, ...)
```

```
clmm_imp(fixed, data, random, n.chains = 3, n.adapt = 100,  
  n.iter = 0, thin = 1, monitor_params = NULL, auxvars = NULL,  
  refcats = NULL, models = NULL, no_model = NULL, trunc = NULL,  
  ridge = FALSE, ppc = TRUE, seed = NULL, inits = NULL,  
  parallel = FALSE, ncores = NULL, scale_vars = NULL,  
  scale_pars = NULL, hyperpars = NULL, modelname = NULL,  
  modeldir = NULL, keep_model = FALSE, overwrite = NULL,  
  quiet = TRUE, progress.bar = "text", warn = TRUE, mess = TRUE,  
  keep_scaled_mcmc = FALSE, ...)
```

```
survreg_imp(formula, data, n.chains = 3, n.adapt = 100, n.iter = 0,  
  thin = 1, monitor_params = NULL, auxvars = NULL, refcats = NULL,  
  models = NULL, no_model = NULL, trunc = NULL, ridge = FALSE,  
  ppc = TRUE, seed = NULL, inits = NULL, parallel = FALSE,  
  ncores = NULL, scale_vars = NULL, scale_pars = NULL,  
  hyperpars = NULL, modelname = NULL, modeldir = NULL,  
  keep_model = FALSE, overwrite = NULL, quiet = TRUE,  
  progress.bar = "text", warn = TRUE, mess = TRUE,  
  keep_scaled_mcmc = FALSE, ...)
```

```
coxph_imp(formula, data, n.chains = 3, n.adapt = 100, n.iter = 0,  
  thin = 1, monitor_params = NULL, auxvars = NULL, refcats = NULL,  
  models = NULL, no_model = NULL, trunc = NULL, ridge = FALSE,  
  ppc = TRUE, seed = NULL, inits = NULL, parallel = FALSE,
```

```
ncores = NULL, scale_vars = NULL, scale_pars = NULL,
hyperpars = NULL, modelname = NULL, modeldir = NULL,
keep_model = FALSE, overwrite = NULL, quiet = TRUE,
progress.bar = "text", warn = TRUE, mess = TRUE,
keep_scaled_mcmc = FALSE, ...)
```

## Arguments

formula	a two sided model formula (see <a href="#">formula</a> )
data	a <code>data.frame</code>
n.chains	the number of parallel chains for the model
n.adapt	the number of iterations for adaptation. See <a href="#">adapt</a> for details. If <code>n.adapt = 0</code> then no adaptation takes place.
n.iter	number of iterations to monitor
thin	thinning interval for monitors
monitor_params	named vector specifying which parameters should be monitored
auxvars	optional vector of variable names that should be used as predictors in the imputation procedure (and will be imputed if necessary) but are not part of the analysis model
refcats	optional; a named list specifying which category should be used as reference category for each of the categorical variables. Options are the category label, the category number, "first" (the first category), "last" (the last category) or "largest" (chooses the category with the most observations). Default is "first". (See also <a href="#">set_refcat</a> )
models	optional named vector specifying the order and types of the models for incomplete covariates and longitudinal covariates. This argument replaces the argument <code>meth</code> used in earlier versions. If <code>NULL</code> (default) models will be determined automatically based on the class of the respective columns of <code>data</code> (see <a href="#">Details</a> ). The default order is incomplete baseline covariates, complete longitudinal covariates, incomplete longitudinal covariates, and within each group variables are ordered according to the proportion of missing values (increasing).
no_model	names of variables for which no model should be specified. Note that this is only possible for completely observed variables and may imply assumptions of independence between the excluded variable and incomplete variables.
trunc	optional named list specifying the limits of truncation for the distribution of the named incomplete variables
ridge	logical; should the parameters of the main model be penalized using ridge regression? Default is <code>FALSE</code>
ppc	logical: should monitors for posterior predictive checks be set? (not yet used)
seed	optional seed value for reproducibility
inits	optional specification of initial values in the form of a list or a function (see <a href="#">jags.model</a> ). If omitted, initial values will be generated automatically by JAGS. It is an error to supply an initial value for an observed node.

parallel	logical; should the chains be sampled using parallel computation? Default is FALSE
ncores	number of cores to use for parallel computation; if left empty all except two cores will be used
scale_vars	optional; named vector of (continuous) variables that will be scaled (such that mean = 0 and sd = 1) to improve convergence of the MCMC sampling. Default is that all continuous variables that are not transformed by a function (e.g. <code>log()</code> , <code>ns()</code> ) will be scaled. Variables for which "lognorm" is set as imputation model are only scaled with regards to the standard deviation, but not centered. Variables to be imputed with "gamma", "beta" or "glmm_gamma" are not scaled. If set to FALSE no scaling will be done.
scale_pars	optional matrix of parameters used for centering and scaling continuous covariates. If not specified, this will be calculated automatically. If FALSE, no scaling will be done.
hyperpars	list of hyperparameters, as obtained by <code>default_hyperpars()</code> ; only needs to be supplied if hyperparameters other than the default should be used
modelName	optional; character string specifying the name of the model file (including the ending, either .R or .txt). If unspecified a random name will be generated.
modelDir	optional; directory containing model file or directory in which the model file will be written. If unspecified a temporary directory will be created.
keep_model	logical; whether the created JAGS model should be saved or removed from the disk (FALSE; default) when the sampling has finished.
overwrite	logical; whether an existing model file with the specified <code>&lt;modelDir&gt;/&lt;modelName&gt;</code> should be overwritten. If set to FALSE and a model already exists, that model will be used. If unspecified (NULL) and a file exists, the user is asked for input on how to proceed
quiet	if TRUE then messages generated during compilation will be suppressed, as well as the progress bar during adaptation.
progress.bar	character string specifying the type of progress bar. Possible values are "text", "gui", and "none". See <a href="#">update</a> .
warn	logical; should warnings be given? Default is TRUE. Note: this applies only to warnings given directly by <b>JointAI</b> .
mess	logical; should messages be given? Default is TRUE. Note: this applies only to messages given directly by <b>JointAI</b> .
keep_scaled_mcmc	should the "original" MCMC sample (i.e., the scaled version returned by <code>coda.samples()</code> ) be kept? (The MCMC sample that is re-scaled to the scale of the data is always kept.)
...	additional, optional arguments
family	only for <code>glm_imp</code> and <code>glmm_imp</code> : a description of the distribution and link function to be used in the model. This can be a character string naming a family function, a family function or the result of a call to a family function. (See <a href="#">family</a> and the 'Details' section below.)
fixed	a two sided formula describing the fixed-effects part of the model (see <a href="#">formula</a> )

random only for lme\_imp, glme\_imp and clmm\_imp: a one-sided formula of the form  $\sim x_1 + \dots + x_n \mid g$ , where  $x_1 + \dots + x_n$  specifies the model for the random effects and  $g$  the grouping variable

### Value

An object of class [JointAI](#).

### Details

See also the vignette: [Model Specification](#)

#### Implemented distribution families and link functions for glm\_imp() and glme\_imp():

gaussian	with links: identity, log
binomial	with links: logit, probit, log, cloglog
Gamma	with links: identity, log
poisson	with links: log, identity

**Imputation methods:** Implemented imputation models that can be chosen in the argument models are:

norm	linear model
lognorm	log-linear model for skewed continuous data
gamma	gamma model (with log-link) for skewed continuous data
beta	beta model (with logit-link) for skewed continuous data in (0, 1)
logit	logistic model for binary data
multilogit	multinomial logit model for unordered categorical variables
cumlogit	cumulative logit model for ordered categorical variables
lmm	linear mixed model for continuous longitudinal covariates
glmm_gamma	Gamma mixed model for skewed longitudinal covariates
glmm_logit	logit mixed model for binary longitudinal covariates
glmm_poisson	Poisson mixed model for longitudinal count covariates
clmm	cumulative logit mixed model for longitudinal ordered factors

When models are specified for only a subset of the incomplete or longitudinal covariates involved in a model, the default choices are used for all unspecified variables.

**Parameters to follow** (monitor\_params): See also the vignette: [Parameter Selection](#)

Named vector specifying which parameters should be monitored. This can be done either directly by specifying the name of the parameter or indirectly by one of the key words selecting a set of parameters. Except for other, in which parameter names are specified directly, parameter (groups) are just set as TRUE or FALSE. If left unspecified, `monitor_params = c("analysis_main" = TRUE)` will be used.

name/key word	what is monitored
analysis_main	betas and sigma_y (and D in mixed models)
analysis_random	ranef, D, invD, RinvD
imp_pars	alphas, tau_imp, gamma_imp, delta_imp

imps	imputed values
betas	regression coefficients of the analysis model
tau_y	precision of the residuals from the analysis model
sigma_y	standard deviation of the residuals from the analysis model
ranef	random effects b
D	covariance matrix of the random effects
invD	inverse of D
RinvD	matrix in the prior for invD
alphas	regression coefficients in the imputation models
tau_imp	precision parameters of the residuals from imputation models
gamma_imp	intercepts in ordinal imputation models
delta_imp	increments of ordinal intercepts
other	additional parameters

For example:

`monitor_params = c(analysis_main = TRUE, tau_y = TRUE, sigma_y = FALSE)` would monitor the regression parameters `betas` and the residual precision `tau_y` instead of the residual standard deviation `sigma_y`.

`monitor_params = c(imps = TRUE)` would monitor `betas`, `tau_y`, and `sigma_y` (because `analysis_main = TRUE` by default) as well as the imputed values.

## Note

**Coding of variables::** The default imputation methods are chosen based on the class of each of the incomplete variables, distinguishing between `numeric`, `factor` with two levels, `unordered factor` with >2 levels and `ordered factor` with >2 levels.

When a continuous incomplete variable has only two different values it is assumed to be binary and its coding and default imputation model will be changed accordingly. This behavior can be overwritten when the imputation method for that variable is specified directly by the user.

Variables of type `'logical'` are automatically be converted to `unordered factors`.

Contrary to base R behavior, dummy coding (i.e., `contr.treatment` contrasts) are used for ordered factors in any linear predictor. However, since the order of levels in an ordered factor contains information relevant to the imputation of missing values, it is important that incomplete ordinal variables are coded as such.

**Non-linear effects and transformation of variables::** **JointAI** handles non-linear effects, transformation of covariates and interactions the following way: When, for instance, the model formula contains the function `log(x)` and `x` has missing values, `x` will be imputed and used in the linear predictor of models for other incomplete variables, i.e., it is assumed that the other variables have a linear association with `x` but not with `log(x)`. The `log()` of the observed and imputed values of `x` is calculated and used in the linear predictor of the analysis model.

If, instead of using `log(x)` in the model formula a pre-calculated variable `logx` would be used instead, this variable is imputed directly and used in the linear predictors of all models, implying that other incomplete variables that have `logx` in their linear predictors have a linear association with `logx` but not with `x`.

When different transformations of the same incomplete variable are used in one model it is strongly discouraged to calculate these transformations beforehand and supply them as different variables. If, for example, a model formula contains both `x` and `x2` (where `x2 = x^2`), they



are treated as separate variables and imputed with separate models. Imputed values of  $x^2$  are thus not equal to the square of imputed values of  $x$ . Instead,  $x + I(x^2)$  should be used in the model formula. Then, only  $x$  is imputed and used in the linear predictor of models for other incomplete variables, and  $x^2$  is calculated from the imputed values of  $x$  internally.

The same applies to interactions involving incomplete variables.

#### Not (yet) possible::

- multiple nesting levels of random effects (nested or crossed)
- prediction (using `predict`) conditional on random effects
- the use of splines for incomplete variables
- the use of `pspline`, `frailty`, `cluster` or `strata` in survival models
- left censored or interval censored data

#### See Also

[set\\_refcat](#), [get\\_models](#), [traceplot](#), [densplot](#), [summary.JointAI](#), [MC\\_error](#), [GR\\_crit](#), [jags.model](#), [coda.samples](#), [predict.JointAI](#), [JointAIObject](#), [add\\_samples](#), [parameters](#), [list\\_impmodels](#)

#### Vignettes

- [Minimal Example](#)
- [Model Specification](#)
- [Parameter Selection](#)
- [After Fitting](#)

#### Examples

```
# Example 1: Linear regression with incomplete covariates
mod1 <- lm_imp(y~C1 + C2 + M2, data = wideDF, n.iter = 100)

# Example 2: Logistic regression with incomplete covariates
mod2 <- glm_imp(B1 ~ C1 + C2 + M2, data = wideDF,
               family = binomial(link = "logit"), n.iter = 100)

# Example 3: Linear mixed model with incomplete covariates
mod3 <- lme_imp(y ~ C1 + B2 + c1 + time, random = ~ time|id,
               data = longDF, n.iter = 500)
```

---

NHANES

*National Health and Nutrition Examination Survey (NHANES) Data*

---

### Description

This data is a small subset of the data collected within the 2011-2012 wave of the NHANES study, a study designed to assess the health and nutritional status of adults and children in the United States, conducted by the [National Center for Health Statistics](#).

### Usage

`data(NHANES)`

### Format

A data frame with 186 rows and 13 variables:

**SBP** systolic blood pressure

**gender** male or female

**age** in years

**race** race / Hispanic origin (5 categories)

**WC** waist circumference in cm

**alc** alcohol consumption (binary: <1 drink per week vs. >= 1 drink per week)

**educ** educational level (binary: low vs. high)

**creat** creatinine concentration in mg/dL

**albu** albumin concentration in g/dL

**uricacid** uric acid concentration in mg/dL

**bili** bilirubin concentration in mg/dL

**occup** occupational status (3 categories)

**smoke** smoking status (3 ordered categories)

### Note

The subset provided here was selected and re-coded to facilitate demonstration of the functionality of the JointAI package, and no clinical conclusions should be derived from it.

### Source

National Center for Health Statistics (NCHS) (2011 - 2012). National Health and Nutrition Examination Survey Data. URL <https://www.cdc.gov/nchs/nhanes/>.

### Examples

`summary(NHANES)`

---

parameters	<i>Parameter names of an JointAI object</i>
------------	---

---

### Description

Returns the names of the parameters/nodes of an object of class 'JointAI' for which a monitor is set.

### Usage

```
parameters(object, mess = TRUE, warn = TRUE)
```

### Arguments

object	object inheriting from class 'JointAI'
mess	logical; should messages be given? Default is TRUE. Note: this applies only to messages given directly by <b>JointAI</b> .
warn	logical; should warnings be given? Default is TRUE. Note: this applies only to warnings given directly by <b>JointAI</b> .
...	currently not used

### Examples

```
# (does not need MCMC samples to work, so we will set n.adapt = 0 and
# n.iter = 0 to reduce computational time)
mod1 <- lm_imp(y ~ C1 + C2 + M2 + O2 + B2, data = wideDF, n.adapt = 0, n.iter = 0)

parameters(mod1)
```

---

plot_all	<i>Visualize the distribution of all variables in the dataset</i>
----------	---

---

### Description

Plots a grid of histograms (for continuous variables) and barplots (for categorical variables) together with the proportion of missing values in each variable.

### Usage

```
plot_all(data, nrow = NULL, ncol = NULL, fill = grDevices::grey(0.8),
  border = "black", allNA = FALSE, use_level = FALSE, idvar,
  xlab = "", ylab = "frequency", ...)
```

**Arguments**

data	a data.frame (or a matrix)
nrow	optional number of rows and columns in the plot layout; automatically chosen if unspecified
ncol	optional number of rows and columns in the plot layout; automatically chosen if unspecified
fill	color the histograms and bars are filled with
border	color of the borders of the histograms and bars
allNA	logical; if FALSE (default) the proportion of missing values is only given for variables that have missing values, if TRUE it is given for all variables
use_level	logical; should the multi-level structure be taken into account? This requires specification of the argument idvar.
idvar	name of the column that specifies the multi-level grouping structure
xlab	labels for the x- and y-axis
ylab	labels for the x- and y-axis
...	additional parameters passed to <code>barplot</code> and <code>hist</code>

**See Also**

Vignette: [Visualizing Incomplete Data](#)

**Examples**

```
par(mar = c(1,2,3,1), mgp = c(2, 0.6, 0))
plot_all(wideDF)
```

---

plot_imp_distr	<i>Plot the distribution of observed and imputed values</i>
----------------	---

---

**Description**

Plots densities and barplots of the observed and imputed values in a long-format dataset (multiple imputed datasets stacked onto each other).#'

**Usage**

```
plot_imp_distr(data, imp = "Imputation_", id = ".id",
  rownr = ".rownr", ncol = NULL, nrow = NULL)
```

**Arguments**

data	a data.frame containing multiple imputations (and the original incomplete data)
imp	the name of the variable specifying the imputation indicator
id	the name of the variable specifying the subject indicator
rownr	the name of a variable identifying which rows correspond to the same observation in the original (unimputed) data
ncol	optional number of rows and columns in the plot layout; automatically chosen if unspecified
nrow	optional number of rows and columns in the plot layout; automatically chosen if unspecified

**Examples**

```
mod <- lme_imp(y ~ C1 + c2 + B2 + b2 + C2 + p2, random = ~ 1 | id, data = longDF,
              n.iter = 500, monitor_params = c(imps = TRUE), mess = FALSE)
impDF <- get_MIdat(mod, m = 5)
plot_imp_distr(impDF, id = "id")
```

---

 predDF

---

*Create a new dataframe for prediction*


---

**Description**

Build a data.frame for prediction, where one variable varies and all other variables are set to the reference value (median for continuous variables).

**Usage**

```
predDF(...)  
  
## S3 method for class 'JointAI'  
predDF(object, var, ...)  
  
## S3 method for class 'formula'  
predDF(formula, dat, var, ...)
```

**Arguments**

...	optional, additional arguments (currently not used)
object	object inheriting from class 'JointAI'
var	name of variable that should be varying
formula	a two sided model formula (see <a href="#">formula</a> )
dat	original data

**See Also**

[predict.JointAI](#), [lme\\_imp](#), [glm\\_imp](#), [lm\\_imp](#)

**Examples**

```
# fit a JointAI model
mod <- lm_imp(y~C1 + C2 + M2, data = wideDF, n.iter = 100)

# generate a dataframe with varying "C2" and reference values for all other variables in the model
newDF <- predDF(mod, var = "C2")

head(newDF)
```

---

<code>predict.JointAI</code>	<i>Predict values from an object of class JointAI</i>
------------------------------	---

---

**Description**

Calculates the expected outcome value for a given set of covariate values and an object of class 'JointAI', and corresponding 2.5% and 97.5% (or other quantiles) credible intervals.

**Usage**

```
## S3 method for class 'JointAI'
predict(object, newdata, quantiles = c(0.025, 0.975),
        start = NULL, end = NULL, thin = NULL, ...)
```

**Arguments**

object	object inheriting from class 'JointAI'
newdata	new dataset for prediction
quantiles	quantiles of the predicted distribution of the outcome
start	the first iteration of interest (see <a href="#">window.mcmc</a> )
end	the last iteration of interest (see <a href="#">window.mcmc</a> )
thin	thinning interval (see <a href="#">window.mcmc</a> )
...	currently not used

**Details**

A model matrix  $X$  is created from the model formula (fixed effects only) and newdata.  $X\beta$  is then calculated for each iteration of the MCMC sample in object, i.e.,  $X\beta$  has `n.iter` rows and `nrow(newdata)` columns. A subset of the MCMC sample can be selected using `start`, `end` and `thin`.

**Value**

A list with entries "fit" and "quantiles", where "fit" contains the column means of  $X\beta$  (see details) and "quantiles" contain the specified quantiles (by default 2.5% and 97.5%) of each column of  $X\beta$ .

**Note**

- For repeated measures models prediction is performed on fixed effects only.
- Prediction is performed on the scale of the linear predictor.

Functionality will be extended in the future.

**See Also**

[predDF.JointAI](#), [lme\\_imp](#), [glm\\_imp](#), [lm\\_imp](#)

**Examples**

```
# fit model
mod <- lm_imp(y ~ C1 + C2 + I(C2^2), data = wideDF, n.iter = 100)

# create dataset for prediction
newDF <- predDF(mod, var = "C2")

# obtain predicted values
pred <- predict(mod, newdata = newDF)

# plot predicted values and 95% confidence band
plot(newDF$C2, pred$fit, type = "l", ylim = range(pred$quantiles),
      xlab = "C2", ylab = "predicted values")
matplot(newDF$C2, t(pred$quantiles), lty = 2, add = TRUE, type = "l", col = 1)
```

---

set\_refcat

*Set the reference categories for all categorical covariates in the model*

---

**Description**

The function asks questions and, depending on the answers given by the user, returns the input for the argument refcats in the functions [lm\\_imp](#), [glm\\_imp](#), [clm\\_imp](#), [lme\\_imp](#), [glme\\_imp](#), [clmm\\_imp](#), [survreg\\_imp](#) and [coxph\\_imp](#).

**Usage**

```
set_refcat(data, formula, covars, auxvars)
```

**Arguments**

data	a data.frame
formula	optional; model formula (used to select subset of relevant columns of data)
covars	optional; vector containing the names of relevant columns of data
auxvars	optional; vector containing the names of relevant columns of data that should be considered additionally to the columns occurring in the formula

**Examples**

```
## Not run:
# Example 1: set reference categories for the whole dataset and choose answer option 3:
set_refcat(data = NHANES)
3

# insert the returned string as argument refcats
mod1 <- lm_imp(SBP ~ age + race + creat + educ, data = NHANES, refcats = 'largest')

# Example 2:
# specify a model formula
fmla <- SBP ~ age + gender + race + bili + smoke + alc

# write the output of set_refcat to an object
ref_mod2 <- set_refcat(data = NHANES, formula = fmla)
4
2
5
1
1

# enter the output in the model specification
mod2 <- lm_imp(formula = fmla, data = NHANES, refcats = ref_mod2, n.adapt = 0)

## End(Not run)
```

---

sharedParams

*Parameters used by several functions in JointAI.*


---

**Description**

Parameters used by several functions in JointAI.

**Arguments**

object	object inheriting from class 'JointAI'
no_model	names of variables for which no model should be specified. Note that this is only possible for completely observed variables and may imply assumptions of independence between the excluded variable and incomplete variables.



subset	subset of parameters/variables/nodes (columns in the MCMC sample). Uses the same logic as the argument <code>monitor_params</code> in <code>lm_imp</code> , <code>glm_imp</code> , <code>clm_imp</code> , <code>lme_imp</code> , <code>glme_imp</code> , <code>survreg_imp</code> and <code>coxph_imp</code> .
start	the first iteration of interest (see <code>window.mcmc</code> )
end	the last iteration of interest (see <code>window.mcmc</code> )
thin	thinning interval (see <code>window.mcmc</code> )
nrow, ncol	optional number of rows and columns in the plot layout; automatically chosen if unspecified
use_ggplot	logical; Should ggplot be used instead of the base graphics?
warn	logical; should warnings be given? Default is TRUE. Note: this applies only to warnings given directly by <b>JointAI</b> .
mess	logical; should messages be given? Default is TRUE. Note: this applies only to messages given directly by <b>JointAI</b> .
xlab, ylab	labels for the x- and y-axis
use_level	logical; should the multi-level structure be taken into account? This requires specification of the argument <code>idvar</code> .
idvar	name of the column that specifies the multi-level grouping structure
keep_aux	logical; Should constant effects of auxiliary variables be kept in the output?
ridge	logical; should the parameters of the main model be penalized using ridge regression? Default is FALSE
parallel	logical; should the chains be sampled using parallel computation? Default is FALSE
ncores	number of cores to use for parallel computation; if left empty all except two cores will be used
seed	optional seed value for reproducibility
ppc	logical: should monitors for posterior predictive checks be set? (not yet used)

---

 simLong

---

*Simulated Longitudinal Data in Long and Wide Format*


---

## Description

This data was simulated to mimic data from a longitudinal cohort study following mothers and their child from birth until approximately 4 years of age. It contains 3903 observations of 499 mother-child pairs. Children's BMI and head circumference was measured repeatedly and their age in months was recorded at each measurement. Furthermore, the data contain several baseline variables with information on the mothers' demographics and socioeconomic status.

## Usage

simLong

simWide

**Format**

simLong: A data frame in long format with 3894 rows and 13 variables

simWide: A data frame in wide format with 499 rows and 45 variables

**Baseline covariates**

(in simLong and simWide)

**GESTBIR** gestational age at birth (in weeks)

**ETHN** ethnicity (binary: European vs. other)

**AGE\_M** age of the mother at intake

**HEIGHT\_M** height of the mother (in cm)

**PARITY** number of times the mother has given birth (binary: 0 vs. >=1)

**SMOKE** smoking status of the mother during pregnancy (3 ordered categories: never smoked during pregnancy, smoked until pregnancy was known, continued smoking in pregnancy)

**EDUC** educational level of the mother (3 ordered categories: low, mid, high)

**MARITAL** marital status (3 categories)

**ID** subject identifier

**Long-format variables**

(only in simLong)

**time** measurement occasion/visit (by design, children should be measured at/around 1, 2, 3, 4, 7, 11, 15, 20, 26, 32, 40 and 50 months of age)

**age** child age at measurement time in months

**bmi** child BMI

**hc** child head circumference in cm

**Wide-format variables**

(only in simWide)

**age1, age2, age3, age4, age7, age11, age15, age20, age26, age32, age40, age50** child age at the repeated measurements in months

**bmi1, bmi2, bmi3, bmi4, bmi7, bmi11, bmi15, bmi20, bmi26, bmi32, bmi40, bmi50** repeated measurements of child BMI

**age1, age2, age3, age4, age7, age11, age15, age20, age26, age32, age40, age50** repeated measurements of child head circumference in cm

**Examples**

```
summary(simLong)
```

```
summary(simWide)
```

---

summary.JointAI      *Summary of an object of class JointAI*

---

## Description

Obtain and print the summary, (fixed effects) coefficients (coef) and credible interval (confint) for an object of class 'JointAI'.

## Usage

```
## S3 method for class 'JointAI'
summary(object, start = NULL, end = NULL,
  thin = NULL, quantiles = c(0.025, 0.975), subset = NULL,
  warn = TRUE, mess = TRUE, ...)

## S3 method for class 'summary.JointAI'
print(x, digits = max(3, .Options$digits - 4),
  ...)

## S3 method for class 'JointAI'
coef(object, start = NULL, end = NULL, thin = NULL,
  subset = NULL, warn = TRUE, mess = TRUE, ...)

## S3 method for class 'JointAI'
confint(object, parm = NULL, level = 0.95,
  quantiles = NULL, start = NULL, end = NULL, thin = NULL,
  subset = NULL, warn = TRUE, mess = TRUE, ...)

## S3 method for class 'JointAI'
print(x, digits = max(4, getOption("digits") - 4), ...)
```

## Arguments

object	object inheriting from class 'JointAI'
start	the first iteration of interest (see <a href="#">window.mcmc</a> )
end	the last iteration of interest (see <a href="#">window.mcmc</a> )
thin	thinning interval (see <a href="#">window.mcmc</a> )
quantiles	posterior quantiles
subset	subset of parameters/variables/nodes (columns in the MCMC sample). Uses the same logic as the argument monitor_params in <a href="#">lm_imp</a> , <a href="#">glm_imp</a> , <a href="#">clm_imp</a> , <a href="#">lme_imp</a> , <a href="#">glme_imp</a> , <a href="#">survreg_imp</a> and <a href="#">coxph_imp</a> .
warn	logical; should warnings be given? Default is TRUE. Note: this applies only to warnings given directly by <b>JointAI</b> .
mess	logical; should messages be given? Default is TRUE. Note: this applies only to messages given directly by <b>JointAI</b> .

...	currently not used
x	an object of class <code>summary.JointAI</code> or <code>JointAI</code>
digits	minimal number of <i>significant</i> digits, see <code>print.default</code> .
parm	same as <code>subset</code>
level	confidence level (default is 0.95)

### See Also

The model fitting functions `lm_imp`, `glm_imp`, `clm_imp`, `lme_imp`, `glme_imp`, `survreg_imp` and `coxph_imp`, and the vignette [Parameter Selection](#) for examples how to specify the parameter subset.

### Examples

```
mod1 <- lm_imp(y~C1 + C2 + M2, data = wideDF, n.iter = 100)

summary(mod1)
coef(mod1)
confint(mod1)
```

---

traceplot

*Traceplot of a JointAI model*

---

### Description

Creates a set of traceplots from the MCMC sample of an object of class "JointAI".

### Usage

```
traceplot(object, ...)

## S3 method for class 'mcmc.list'
traceplot(object, start = NULL, end = NULL,
  thin = NULL, ...)

## S3 method for class 'JointAI'
traceplot(object, start = NULL, end = NULL,
  thin = NULL, subset = c(analysis_main = TRUE), nrow = NULL,
  ncol = NULL, keep_aux = FALSE, use_ggplot = FALSE, warn = TRUE,
  ...)
```

**Arguments**

object	object inheriting from class 'JointAI'
...	Arguments passed on to <code>graphics::matplot</code>
	<b>lty</b> vector of line types, widths, and end styles. The first element is for the first column, the second element for the second column, etc., even if lines are not plotted for all columns. Line types will be used cyclically until all plots are drawn.
	<b>lwd</b> vector of line types, widths, and end styles. The first element is for the first column, the second element for the second column, etc., even if lines are not plotted for all columns. Line types will be used cyclically until all plots are drawn.
	<b>lend</b> vector of line types, widths, and end styles. The first element is for the first column, the second element for the second column, etc., even if lines are not plotted for all columns. Line types will be used cyclically until all plots are drawn.
	<b>col</b> vector of colors. Colors are used cyclically.
	<b>cex</b> vector of character expansion sizes, used cyclically. This works as a multiple of <code>par("cex")</code> . NULL is equivalent to 1.0.
	<b>bg</b> vector of background (fill) colors for the open plot symbols given by <code>pch = 21:25</code> as in <code>points</code> . The default NA corresponds to the one of the underlying function <code>plot.xy</code> .
	<b>xlim</b> ranges of x and y axes, as in <code>plot</code> .
	<b>ylim</b> ranges of x and y axes, as in <code>plot</code> .
	<b>add</b> logical. If TRUE, plots are added to current one, using <code>points</code> and <code>lines</code> .
	<b>verbose</b> logical. If TRUE, write one line of what is done.
start	the first iteration of interest (see <code>window.mcmc</code> )
end	the last iteration of interest (see <code>window.mcmc</code> )
thin	thinning interval (see <code>window.mcmc</code> )
subset	subset of parameters/variables/nodes (columns in the MCMC sample). Uses the same logic as the argument <code>monitor_params</code> in <code>lm_imp</code> , <code>glm_imp</code> , <code>clm_imp</code> , <code>lme_imp</code> , <code>glme_imp</code> , <code>survreg_imp</code> and <code>coxph_imp</code> .
nrow	optional number of rows and columns in the plot layout; automatically chosen if unspecified
ncol	optional number of rows and columns in the plot layout; automatically chosen if unspecified
keep_aux	logical; Should constant effects of auxiliary variables be kept in the output?
use_ggplot	logical; Should ggplot be used instead of the base graphics?
warn	logical; should warnings be given? Default is TRUE. Note: this applies only to warnings given directly by <b>JointAI</b> .

**See Also**

`summary.JointAI`, `lme_imp`, `glm_imp`, `lm_imp`, `densplot` The vignette **Parameter Selection** contains some examples how to specify the parameter subset.

**Examples**

```
# fit a JointAI model
mod <- lm_imp(y~C1 + C2 + M2, data = wideDF, n.iter = 100)

# Example 1: simple traceplot
traceplot(mod)

# Example 2: ggplot version of traceplot
traceplot(mod, use_ggplot = TRUE)

# Example 5: changing how the ggplot version looks (using standard ggplot syntax)
library(ggplot2)

traceplot(mod, use_ggplot = TRUE) +
  theme(legend.position = 'botto') +
  xlab('iteration') +
  ylab('value') +
  scale_color_discrete(name = 'chain')
```

---

wideDF

*Cross-sectional example dataset*


---

**Description**

A simulated cross-sectional dataset.

**Usage**

```
data(wideDF)
```

**Format**

A simulated data frame with 100 rows and 13 variables:

- C1** continuous, complete variable
- C2** continuous, incomplete variable
- B1** binary, complete variable
- B2** binary, incomplete variable
- M1** unordered factor; complete variable
- M2** unordered factor; incomplete variable
- O1** ordered factor; complete variable
- O2** ordered factor; incomplete variable

- L1** continuous, complete variable
- L2** continuous incomplete variable
- id** id (grouping) variable
- time** continuous complete variable
- y** continuous, complete variable

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