Package ‘mitml’

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**mitml-package**

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**Description**

Provides tools for multiple imputation of missing data in multilevel modeling. This package includes a user-friendly interface to the algorithms implemented in the R packages pan and jomo, as well as several functions for visualizing, managing and analyzing multiply imputed data sets.

The main interface to pan is the function `panImpute`, which allows specification of imputation models for continuous variables with missing data at level 1. In addition, the function `jomoImpute` provides an interface to jomo, which extends the functionality of pan to continuous and categorical variables with missing data at level 1 and level 2 and also allows for the specification of single-level imputation models. Imputations and parameter chains are stored in objects of class `mitml`. To obtain the completed (i.e., imputed) data sets, `mitmlComplete` is used, producing a list of imputed data sets of class `mitmlNlist` that can be used in further analyses.

Several additional functions allow for convenient analysis of multiply imputed data sets, especially when using the R packages `lme4` and `nlme`. The functions `within`, `sort`, and `subset` can be used to manage and manipulate multiply imputed data sets. For model fitting, `with` is used. Final parameter estimates can be extracted using `testEstimates`. Single- and multi-parameter hypotheses tests can be performed using the functions `testConstraints` and `testModels`. In addition, the `anova`
method provides a simple interface to model comparisons with automatic refitting of statistical models.

Data sets can be imported and exported from or to different statistical software packages. Currently, mids2mitml.list, amelia2mitml.list, jomo2mitml.list, and long2mitml.list can be used for importing imputations for other packages in R. In addition, write.mitmlMplus, write.mitmlSAV, and write.mitmlSPSS export data sets to Mplus and SPSS, respectively.

Finally, the package provides tools for summarizing and visualizing imputation models, which is useful for the assessment of convergence and the reporting of results.

The data sets contained in this package are published under the same license as the package itself. They contain simulated data and may be used by anyone free of charge as long as reference to this package is given.

Author(s)

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Maintainer: Simon Grund <grund@ipn.uni-kiel.de>

amelia2mitml.list

Convert objects of class amelia to mitml.list

Description

This function converts a amelia class object (as produced by the Amelia package) to mitml.list. The resulting object may be used in further analyses.

Usage

amelia2mitml.list(x)

Arguments

x An object of class amelia as produced by amelia (see the Amelia package).

Value

A list of imputed data sets with an additional class attribute mitml.list.

Author(s)

Simon Grund

See Also

mitmlComplete
Examples

```r
data(studentratings)
require(Amelia)
imp <- amelia(x=studentratings[,c("ID","MathAchiev","ReadAchiev")], cs="ID")
implist <- amelia2mitml.list(imp)
```

Description

Performs model comparisons for a series of nested statistical models fitted using `with.mitml.list`.

Usage

```r
# S3 method for class 'mitml.result'
anova(object, ...)
```

Arguments

- `object` An object of class `mitml.result` as produced by `with.mitml.list`.
- `...` Additional objects of class `mitml.result` to be included in the comparison.

Details

This function performs several model comparisons between models fitted using `with.mitml.list`. If possible, the models are compared using the $D_3$ statistic (Meng & Rubin, 1992). If this method is unavailable, the $D_2$ statistic is used instead (Li, Meng, Raghunathan, & Rubin, 1991). The $D_3$ method currently supports linear models and linear mixed-effects models with a single cluster variable as estimated by `lme4` or `nlme` (see Laird, Lange, & Stram, 1987).

This function is essentially a wrapper for `testModels` with the advantage that several models can be compared simultaneously. All model comparisons are likelihood-based. For further options for model comparisons (e.g., Wald-based procedures) and finer control, see `testModels`.

Value

Returns a list containing the results of each model comparison. A `print` method is used for better readable console output.

Author(s)

Simon Grund
References


See Also

`withNmitmlNlist`, `testModels`

Examples

```r
require(lme4)
data(studentratings)

fml <- ReadDis + SES ~ ReadAchiev + (1|ID)
imp <- panImpute(studentratings, formula=fml, n.burn=1000, n.iter=100, m=5)

implist <- mitmlComplete(imp, print=1:5)

# simple comparison (same as testModels)
fit0 <- with(implist, lmer(ReadAchiev ~ (1|ID), REML=FALSE))
fit1 <- with(implist, lmer(ReadAchiev ~ ReadDis + (1|ID), REML=FALSE))
anova(fit1,fit0)

## Not run:
# multiple comparisons
fit2 <- with(implist, lmer(ReadAchiev ~ ReadDis + (1+ReadDis|ID), REML=FALSE))
anova(fit2,fit1,fit0)

## End(Not run)
```

---

**as.mitml.list**

*Convert a list of data sets to mitml.list*

**Description**

This function adds a `mitml.list` class attribute to a list of data frames. The resulting object can be used in further analyses.

**Usage**

```r
as.mitml.list(x)
```
Arguments

x

A list of data frames.

Value

The original list with an additional class attribute `mitml.list`. The list entries are converted to `data.frame` if necessary, in which case a note is printed.

Author(s)

Simon Grund

See Also

`is.mitml.list`, `long2mitml.list`

Examples

```r
# data frame with 'imputation' indicator
dat <- data.frame(imputation=rep(1:10,each=20), x=rnorm(200))

# split into a list and convert to 'mitml.list'
l <- split(dat, dat$imputation)
l <- as.mitml.list(l)

is.mitml.list(l)
# TRUE
```

c.mitml.list

---

**Concatenate lists of imputed data sets**

Description

These functions allow concatenating lists of imputed data sets by data set, row, or column.

Usage

```r
## S3 method for class 'mitml.list'
c(...)
## S3 method for class 'mitml.list'
rbind(...)
## S3 method for class 'mitml.list'
cbind(...)
```
Arguments

... One or several lists of imputed data sets with class mitml.list as produced by mitmlComplete (or similar).

Details

These function allow concatenating multiple lists of imputed data sets. The function c concatenates by data set (i.e., by appending additional data sets to the list), rbind concatenates by row (i.e., appending additional rows to each data set), and cbind concatenates by column (i.e., by appending additional columns to each data set).

These functions are intended for experienced users and should be used with caution. Appending rows or columns from multiple imputation procedures is usually unsafe unless in special applications (see Examples).

Value

A list of imputed data sets with an additional class attribute mitml.list.

Author(s)

Simon Grund

Examples

# Example 1: manual imputation by grouping variable
data(studentratings)
fm1 <- ReadDis + SES ~ ReadAchiev + (1|ID)

imp <- panImpute(subset(studentratings, FedState=="SH"), formula=fm1, n.burn=1000, n.iter=100, m=5)
implist <- mitmlComplete(imp, print="all")

imp2 <- panImpute(subset(studentratings, FedState=="B"), formula=fm1, n.burn=1000, n.iter=100, m=5)
implist2 <- mitmlComplete(imp2, print="all")
rbind(implist, implist2)

# Example 2: predicted values from linear model
pred <- with(implist, predict(lm(ReadDis~ReadAchiev)))
cbind(implist, pred.ReadDis=pred)

calculateMeans | Calculate cluster means

Description

Calculates the mean of a given variable within each cluster, possibly conditioning on an additional grouping variable.
clusterMeans

Usage

clusterMeans(x, cluster, adj=FALSE, group=NULL)

Arguments

x  A numeric vector for which cluster means should be calculated. Can also be supplied as a character string denoting a variable in the current environment (see details).

cluster  A numeric vector or a factor denoting the cluster membership of each unit in x. Can also be supplied as a character string (see details).

adj  Logical flag indicating if person-adjusted group means should be calculated. The cluster mean is then calculated for each unit by excluding that unit from calculating the cluster mean. Default is to FALSE.

group  (optional) An grouping factor or a variable that can be interpreted as such. If specified, then cluster means are calculated conditionally on the grouping variable, that is, separately within sub-groups. Can also be supplied as a character string (see details).

Details

This function calculates the mean of a variable within each level of a cluster variable. Any NA are omitted during calculation.

The three main arguments of the function can also be supplied as (single) character strings, denoting the name of the respective variables in the current environment. This is especially useful for calculating several cluster means simultaneously, for example using within.mitml.list (see Example 2 below).

Value

Returns a numeric vector with the same length as x containing the cluster mean for all units.

Author(s)

Simon Grund, Alexander Robitzsch

See Also

within.mitml.list

Examples

data(studentratings)

fml <- ReadDis + SES ~ ReadAchiev + (1|ID)
imp <- panImpute(studentratings, formula=fml, n.burn=1000, n.iter=100, m=5)
implist <- mitmlComplete(imp, print=1:5)

# * Example 1: single cluster means
# calculate cluster means (for each data set)
with(implist, clusterMeans(ReadAchiev, ID))

# ... person-adjusted cluster means
with(implist, clusterMeans(ReadAchiev, ID, adj=TRUE))

# ... groupwise cluster means
with(implist, clusterMeans(ReadAchiev, ID, group=Sex))

# * Example 2: automated cluster means using 'for' and 'assign'
# calculate multiple cluster means within multiply imputed data sets
within(implist,
  vars <- c("ReadAchiev","MathAchiev","CognAbility")
  for(i in vars) assign(paste(i,"Mean",sep="."), clusterMeans(i,ID))
  rm(i,vars)
)

confint.mitml.testEstimates

* Compute confidence intervals *

Description

Computes confidence intervals on the basis of the final parameter estimates and inferences given by *testEstimates*.

Usage

```r
## S3 method for class 'mitml.testEstimates'
confint(object, parm, level=0.95, ...)
```

Arguments

- `object`: An object of class `mitml.testEstimates` as produced by `testEstimates`.
- `parm`: (optional) A reference to the parameters for which to calculate confidence intervals. Can be a character or integer vector denoting names or position of parameters, respectively. If missing, all parameters are considered (the default).
- `level`: The confidence level. Default is to 0.95 (i.e., 95%).
- `...`: Not being used.
Details
This function calculates confidence intervals with the given confidence level for the specified parameters on the basis of a $t$-distribution, with estimates, standard errors, and degrees of freedom as returned by `testestimates`.

Value
A matrix containing the lower and upper bounds of the confidence intervals.

Author(s)
Simon Grund

See Also
`testestimates`

Examples
```r
data(studentratings)

fml <- ReadDis ~ ReadAchiev + (1|ID)
imp <- panImpute(studentratings, formula=fml, n.burn=500, n.iter=100, m=5)
implist <- mitmlComplete(imp, print=1:5)

# fit regression model
fit <- with(implist, lm(ReadDis ~ 1 + ReadAchiev))
est <- testestimates(fit)

# calculate confidence intervals
confint(est)

# ... with different confidence levels
confint(est, level=0.90)
confint(est, level=0.999)
```

---

is.mitml.list

Check if an object is of class `mitml.list`

Description
This function checks if its argument is a list of class `mitml.list`.

Usage

```r
is.mitml.list(x)
```
**jomoImpute**

### Arguments

- **x**  
  An R object.

### Value

Either TRUE or FALSE. A warning message is displayed if the contents of x do not appear to be data frames.

### Author(s)

Simon Grund

### See Also

as.mitml.list

### Examples

```r
1 <- list(data.frame(x=rnorm(20)))
1 <- as.mitml.list(1)
is.mitml.list(1)  
# TRUE

1 <- as.list(1:10)
is.mitml.list(1)  
# FALSE

class(l) <- "mitml.list"
is.mitml.list(l)  
# TRUE, with a warning
```

---

**jomoImpute**

*Impute single-level and multilevel missing data using jomo*

### Description

This function provides an interface to the **jomo** package, which uses the MCMC algorithms presented in Carpenter & Kenward (2013). Through this wrapper function, single-level and multilevel imputations can be generated for (mixed) categorical and continuous variables (Goldstein et al., 2009). The multilevel procedures support imputation of missing data at level 1 and 2 as well as imputation using random (residual) covariance matrices (Yucel, 2011). Imputations can be generated using type or formula, which offer different options for model specification.

### Usage

```r
jomoImpute(data, type, formula, random=L1=c("none","mean","full"), n.burn=5000,  
n.iter=100, m=10, group=NULL, prior=NULL, seed=NULL, save.pred=FALSE,  
keep.chains=c("full","diagonal"), silent=FALSE)
```
Arguments

data  A data frame containing incomplete and auxiliary variables, the cluster indicator variable, and any other variables that should be present in the imputed datasets.

type  An integer vector specifying the role of each variable in the imputation model or a list of two vectors specifying a two-level model (see details).

formula  A formula specifying the role of each variable in the imputation model or a list of two formulas specifying a two-level model. The basic model is constructed by `model.matrix`, which allows including derived variables in the imputation model using `I()` (see details and examples).

random.L1  A character string denoting if the covariance matrix of residuals should vary across groups and how the values of these matrices are stored (see details). Default is to "none", assuming a common covariance matrix across clusters.

n.burn  The number of burn-in iterations before any imputations are drawn. Default is to 5,000.

n.iter  The number of iterations between imputations. Default is to 100.

m  The number of imputed data sets to generate. Default is to 10.

group  (optional) A character string denoting the name of an additional grouping variable to be used with the `formula` argument. When specified, the imputation model is run separately within each of these groups.

prior  (optional) A list with components `binv`, `dinv`, and `a` for specifying prior distributions for the covariance matrix of random effects and the covariance matrix of residuals (see details). Default is to using least-informative priors.

seed  (optional) An integer value initializing R’s random number generator for reproducible results. Default is to using the global seed.

save.pred  (optional) Logical flag indicating if variables derived using `formula` should be included in the imputed data sets. Default is to `FALSE`.

keep.chains  (optional) A character string denoting which parameter chains to save. Default is to save all chains (see details).

silent  (optional) Logical flag indicating if console output should be suppressed. Default is to `FALSE`.

Details

This function serves as an interface to the `jomo` package and supports imputation of single-level and multilevel continuous and categorical data. In order for categorical variables to be detected correctly, these must be formatted as a `factor` variables (see examples). The imputation model can be specified using either the `type` or the `formula` argument.

The `type` interface is designed to provide quick-and-easy imputations using `jomo`. The `type` argument must be an integer vector denoting the role of each variable in the imputation model:

- 1: target variables containing missing data
- 2: predictors with fixed effect on all targets (completely observed)
- 3: predictors with random effect on all targets (completely observed)
- −1: grouping variable within which the imputation is run separately
At least one target variable and, for multilevel imputation, the cluster indicator must be specified. If the cluster indicator is omitted, single-level imputation will be performed. The intercept is automatically included both as a fixed and random effect. If a variable of type \(-1\) is found, then separate imputations are performed within each level of that variable.

The `formula` argument is intended as more flexible and feature-rich interface to `jomo`. Specifying the `formula` argument is similar to specifying other formulae in R. Given below is a list of operators that `jomoImppute` currently understands:

- `~`: separates the target (left-hand) and predictor (right-hand) side of the model
- `+`: adds target or predictor variables to the model
- `*`: adds an interaction term of two or more predictors
- `|`: denotes cluster-specific random effects and specifies the cluster indicator (e.g., `1|ID`)
- `I()`: defines functions to be interpreted by `model.matrix`

If the cluster indicator is omitted, single-level imputation will be run. For multilevel imputation, predictors are allowed to have fixed effects, random effects, or both on all target variables. The intercept is automatically included both as a fixed and, for multilevel imputation, a random effect. Both can be constrained if necessary (see `panImpute`). Note that, when specifying random effects other than the intercept, these will not be automatically added as fixed effects and must be included explicitly. Any predictors defined by `I()` will be used for imputation but not included in the data set unless `save.pred=TRUE`.

If missing data occur at both levels of the sample (level 1 and level 2), then a list of two `formula` s or `type` s may be provided. The first element of this list denotes the imputation model for variables at level 1. The second element denotes the imputation model for variables at level 2. In such a case, missing values are imputed jointly at both levels (see examples, see also Carpenter and Kenward, 2013; Goldstein et al., 2009).

It is possible to model the covariance matrix of residuals at level 1 as random across clusters (Yucel, 2011; Carpenter & Kenward, 2013). The `random.L1` argument determines this behavior and how the values of these matrices are stored. If set to "none", a common covariance matrix is assumed across groups (similar to `panImpute`). If set to "mean", the covariance matrices are random, but only the average covariance matrix is stored at each iteration. If set to "full", the covariance matrices are random, and all variances and covariances from all clusters are stored.

In order to run separate imputations for each level of an additional grouping variable, the `group` argument may be used. The name of the grouping variable must be given in quotes.

As a default prior, `jomoImppute` uses "least informative" inverse-Wishart priors for the covariance matrix of random effects (and residuals at level 2) and the covariance matrix of residuals at level 1, that is, with minimum degrees of freedom (largest dispersion) and identity matrices for scale. For better control, the `prior` argument may be used for specifying alternative prior distributions. These must be supplied as a list containing the following components:

- `Binv`: scale matrix for the covariance matrix of residuals at level 1
- `Dinv`: scale matrix for the covariance matrix of random effects and residuals at level 2
- `a`: starting value for the degrees of freedom of random covariance matrices of residuals (only used with `random.L1="mean"` or `random.L1="full"`)
Note that jomo does not allow for the degrees of freedom for the inverse-Wishart prior to be specified by the user. These are always set to the lowest value possible (largest dispersion) or determined iteratively if the residuals at level 1 are modeled as random (see above). For single-level imputation, only Binv is relevant.

In imputation models with many parameters, the number of parameter chains being saved can be reduced with the keep.chains argument (see panImpute). This setting influences the storage mode of parameters (e.g., dimensions and indices of arrays) and should be used with caution.

Value

Returns an object of class mitml, containing the following components:

data          The original (incomplete) data set, sorted according to the cluster variable and (if given) the grouping variable, with several attributes describing the original order ("sort"), grouping ("group") and factor levels of categorical variables.
replacement.mat A matrix containing the multiple replacements (i.e., imputations) for each missing value. The replacement matrix contains one row for each missing value and one one column for each imputed data set.
index.mat     A matrix containing the row and column index for each missing value. The index matrix is used to link the missing values in the data set with their corresponding rows in the replacement matrix.
call          The matched function call.
model         A list containing the names of the cluster variable, the target variables, and the predictor variables with fixed and random effects, at level 1 and level 2, respectively.
random.L1     A character string denoting the handling of the (random) covariance matrix of residuals at level 1 (see details).
prior         The prior parameters used in the imputation model.
iter          A list containing the number of burn-in iterations, the number of iterations between imputations, and the number of imputed data sets.
par.burnin     A multi-dimensional array containing the parameters of the imputation model from the burn-in phase.
par.imputation A multi-dimensional array containing the parameters of the imputation model from the imputation phase.

Note

For objects of class mitml, methods for the generic functions print, summary, and plot have been defined. mitmlComplete is used for extracting the imputed data sets.

Author(s)

Simon Grund, Alexander Robitzsch, Oliver Luedtke
References


See Also

`panImpute`, `mitmlComplete`, `summary.mitml`, `plot.mitml`

Examples

```r
# NOTE: The number of iterations in these examples is much lower than it
# should be! This is done in order to comply with CRAN policies, and more
# iterations are recommended for applications in practice!

data(studentratings)
data(leadership)

# ***
# for further examples, see "panImpute"
#
?

# *** ........................................
# the 'type' interface
#
#
# * Example 1.1 (studentratings): 'ReadDis' and 'SES', predicted by 'ReadAchiev'
# (random slope)

type <- c(-2,0,0,0,0,1,3,1,0,0)
names(type) <- colnames(studentratings)
type

imp <- jomoImpute(studentratings, type=type, n.burn=100, n.iter=10, m=5)

# * Example 1.2 (leadership): all variables (mixed continuous and categorical
# data with missing values at level 1 and level 2)

type.L1 <- c(-2,1,0,1,1)  # imputation model at level 1
type.L2 <- c(-2,0,1,0,0)  # imputation model at level 2
names(type.L1) <- names(type.L2) <- colnames(leadership)
type <- list(type.L1, type.L2)
type

imp <- jomoImpute(leadership, type=type, n.burn=100, n.iter=10, m=5)
```
Example data set on employees' justice perceptions and satisfaction

**Description**

This data set contains simulated data for employees nested within organizations, featuring employees' sex, ratings on individual justice orientation and ratings on job satisfaction. In addition, the
The data set features scores for justice climate in each organization (defined at the level of organizations, level 2). Different organizations are denoted by the variable id.

The data were simulated based on the results by Liao and Rupp (2005), as well as the secondary analyses of the same data given in Mathieu, Aguinis, Culpepper, and Chen, (2012).

Usage

```r
data(justice)
```

Format

A data frame containing 1400 observations on 4 variables.

References


---

The original data set is available in the `multilevel` package and was altered by (a) transforming workload into a categorical variable, (b) transforming cohesion into a group-level variable, and (c) by inducing missing values.

Usage

```r
data(leadership)
```

Format

A data frame containing 750 observations on 5 variables.
References


---

long2mitml.list

Convert imputations from long format to mitml.list

Description

These functions convert data sets containing multiple imputations in long format to objects of class mitml.list. The resulting object can be used in further analyses.

Usage

long2mitml.list(x, split, exclude=NULL)

jomo2mitml.list(x)

Arguments

x
A data frame in long format containing multiple imputations (see details).

split
A character string denoting the column in x that identifies different imputations (see details).

exclude
A vector denoting values of split which should be excluded from the list.

Details

The function long2mitml.list is intended for converting data frames from the long format to mitml.list (i.e., a list of imputed data sets). In this format, imputations are enclosed in a single data frame, and a split variable is used to identify different imputations.

Similarly, jomo2mitml.list is a special case of long2mitml.list intended for converting imputations that have been generated with jomo directly.

Value

A list of imputed data sets with an additional class attribute mitml.list.

Author(s)

Simon Grund
See Also

mitmlComplete

Examples

data(studentratings)
require(jomo)

# impute data using jomo (native functions)
clus <- studentratings[,"ID"]
Y <- studentratings[,c("ReadAchiev","ReadDis")]
imp <- jomo(Y=Y, clus=clus, nburn=1000, nbetween=100, nimp=5)

# split imputations
implist <- long2mitml.list(imp, split="Imputation", exclude=0)
implist <- jomo2mitml.list(imp)

mids2mitml.list

Convert objects of class mids to mitml.list

Description

This function converts a mids class object (as produced by the mice package) to mitml.list. The resulting object may be used in further analyses.

Usage

mids2mitml.list(x)

Arguments

x An object of class mids as produced by mice (see the mice package).

Value

A list of imputed data sets with an additional class attribute mitml.list.

Author(s)

Simon Grund

See Also

mitmlComplete
Examples

data(studentratings)

# imputation using mice
require(mice)
imp <- mice(studentratings)

implist <- mids2mitml.list(imp)

mitmlComplete   Extract imputed data sets

Description

This function extracts imputed data sets from mitml class objects as produced by panImpute and jomoImpute.

Usage

mitmlComplete(x, print="all", force.list=FALSE)

Arguments

x An object of class mitml as produced by panImpute and jomoImpute.

print Either an integer vector, "list", or "all" denoting which data sets to extract. If set to "list" or "all", then all imputed data sets will be returned as a list. Negative values and zero will return the original (incomplete) data set. Default is to "all".

force.list (optional) Logical flag indicating if single data sets should be enclosed in a list. Default is to FALSE.

Value

Single data sets are returned as a data frame unless force.list=TRUE. If several data sets are extracted, the result is always a list of data sets with an additional class attribute mitml.list.

Author(s)

Simon Grund

See Also

panImpute, jomoImpute
Examples

```r
data(studentratings)

fml <- ReadDis + SES ~ ReadAchiev + (1|ID)
imp <- panImpute(studentratings, formula=fml, n.burn=1000, n.iter=100, m=5)

# extract original (incomplete) data set
mitmComplete(imp, print=0)

# extract first imputed data set (returned as mitml.list)
mitmComplete(imp, print=1, force.list=TRUE)

# extract all imputed data sets at once
implist <- mitmComplete(imp, print="all")

## Not run:
## ... alternatives with same results
implist <- mitmComplete(imp, print=1:5)
implist <- mitmComplete(imp, print="list")

## End(Not run)
```

### multilevelR2

**Calculate R-squared measures for multilevel models**

**Description**

Calculates several measures for the proportion of explained variance in a fitted linear mixed-effects (i.e.,) multilevel model (or a list of fitted models).

**Usage**

```r
multilevelR2(model, print=c("RB1","RB2","SB","MVP"))
```

**Arguments**

- `model`: Either a fitted linear mixed-effects model as produced by `lme4` or `nlme`, or a list of fitted models as produced by `with.mitml.list`.
- `print`: A character vector denoting which measures should be calculated (see details). Default is to printing all measures.

**Details**

This function calculates several measures of explained variance ($R^2$) for linear-mixed effects models. It can be used with a single model, as produced by the packages `lme4` or `nlme`, or a list of
fitted models produced by `with.mitml.list`. In the latter case, the $R^2$ measures are calculated separately for each imputed data set and then averaged across data sets.

Different $R^2$ measures can be requested using the `print` argument. Specifying `RB1` and `RB2` will return the explained variance at level 1 and level 2, respectively, according to Raudenbush and Bryk (2002, pp. 74 and 79). Specifying `SB` will return the total variance explained according to Snijders and Bosker (2012, p. 112). Specifying `MVP` will return the total variance explained based on “multilevel variance partitioning” as proposed by LaHuis, Hartman, Hakoyama, and Clark (2014).

**Value**

Returns a numeric vector containing the $R^2$ measures requested in `print`.

**Note**

Calculating $R^2$ measures is currently only supported for two-level models with a single cluster variable.

**Author(s)**

Simon Grund

**References**


**Examples**

```r
require(lme4)
data(studentratings)

fml <- MathAchiev + ReadAchiev + CognAbility ~ 1 + (1|ID)
imp <- panImpute(studentratings, formula=fml, n.burn=1000, n.iter=100, m=5)
implist <- mitmlComplete(imp, print=1:5)
fit <- with(implist, lmer(MathAchiev ~ 1 + CognAbility + (1|ID)))
multilevelR2(fit)
```
panImpute

**Impute multilevel missing data using pan**

**Description**

This function provides an interface to the pan package for multiple imputation of multilevel data (Schafer & Yucel, 2002). Imputations can be generated using type or formula, which offer different options for model specification.

**Usage**

```r
panImpute(data, type, formula, n.burn=5000, n.iter=100, m=10, group=NULL, prior=NULL, seed=NULL, save.pred=FALSE, keep.chains=c("full","diagonal"), silent=FALSE)
```

**Arguments**

- **data**: A data frame containing incomplete and auxiliary variables, the cluster indicator variable, and any other variables that should be present in the imputed datasets.
- **type**: An integer vector specifying the role of each variable in the imputation model (see details).
- **formula**: A formula specifying the role of each variable in the imputation model. The basic model is constructed by `model.matrix`, thus allowing to include derived variables in the imputation model using `I()` (see details and examples).
- **n.burn**: The number of burn-in iterations before any imputations are drawn. Default is to 5,000.
- **n.iter**: The number of iterations between imputations. Default is to 100.
- **m**: The number of imputed data sets to generate.
- **group**: (optional) A character string denoting the name of an additional grouping variable to be used with the `formula` argument. When specified, the imputation model is run separately within each of these groups.
- **prior**: (optional) A list with components `a`, `Binv`, `c`, and `Dinv` for specifying prior distributions for the covariance matrix of random effects and the covariance matrix of residuals (see details). Default is to using least-informative priors.
- **seed**: (optional) An integer value initializing pan’s random number generator for reproducible results. Default is to using random seeds.
- **save.pred**: (optional) Logical flag indicating if variables derived using `formula` should be included in the imputed data sets. Default is to `FALSE`.
- **keep.chains**: (optional) A character string denoting which parameter chains to save. Default is to save all chains (see details).
- **silent**: (optional) Logical flag indicating if console output should be suppressed. Default is to `FALSE`. 
Details

This function serves as an interface to the `pan` algorithm. The imputation model can be specified using either the `type` or the `formula` argument.

The `type` interface is designed to provide quick-and-easy imputations using `pan`. The `type` argument must be an integer vector denoting the role of each variable in the imputation model:

- 1: target variables containing missing data
- 2: predictors with fixed effect on all targets (completely observed)
- 3: predictors with random effect on all targets (completely observed)
- 4: grouping variable within which the imputation is run separately
- 5: cluster indicator variable
- 0: variables not featured in the model

At least one target variable and the cluster indicator must be specified. The intercept is automatically included both as a fixed and random effect. If a variable of type 4 is found, then separate imputations are performed within each level of that variable.

The `formula` argument is intended as more flexible and feature-rich interface to `pan`. Specifying the `formula` argument is similar to specifying other formulae in R. Given below is a list of operators that `panImpute` currently understands:

- `~`: separates the target (left-hand) and predictor (right-hand) side of the model
- `+`: adds target or predictor variables to the model
- `*`: adds an interaction term of two or more predictors
- `|`: denotes cluster-specific random effects and specifies the cluster indicator (e.g., `1|ID`)
- `I()`: defines functions to be interpreted by `model.matrix`

Predictors are allowed to have fixed effects, random effects, or both on all target variables. The intercept is automatically included both as a fixed and random effect, but it can be constrained if necessary (see examples). Note that, when specifying random effects other than the intercept, these will not be automatically added as fixed effects and must be included explicitly. Any predictors defined by `I()` will be used for imputation but not included in the data set unless `save.pred=TRUE`.

In order to run separate imputations for each level of an additional grouping variable, the `group` argument may be used. The name of the grouping variable must be given in quotes.

As a default prior, `panImpute` uses "least informative" inverse-Wishart priors for the covariance matrix of random effects and the covariance matrix of residuals, that is, with minimum degrees of freedom (largest dispersion) and identity matrices for scale. For better control, the `prior` argument may be used for specifying alternative prior distributions. These must be supplied as a list containing the following components:

- `a`: degrees of freedom for the covariance matrix of residuals
- `Binv`: scale matrix for the covariance matrix of residuals
- `c`: degrees of freedom for the covariance matrix of random effects
- `Dinv`: scale matrix for the covariance matrix of random effects
A sensible choice for a diffuse non-default prior is to set the degrees of freedom to the lowest value possible, and the scale matrices according to a prior guess of the corresponding covariance matrices (see Schafer & Yucel, 2002).

In imputation models with many parameters, the number of parameter chains being saved can be reduced with the `keepNchains` argument. If set to `full` (the default), all chains are saved. If set to `diagonal`, only chains pertaining to fixed effects and the diagonal entries of the covariance matrices are saved. This setting influences the storage mode of parameters (e.g., dimensions and indices of arrays) and should be used with caution.

### Value

Returns an object of class `mitml`, containing the following components:

- **data**: The original (incomplete) data set, sorted according to the cluster variable and (if given) the grouping variable, with several attributes describing the original row order ("sort") and grouping ("group").

- **replacement.mat**: A matrix containing the multiple replacements (i.e., imputations) for each missing value. The replacement matrix contains one row for each missing value and one one column for each imputed data set.

- **index.mat**: A matrix containing the row and column index for each missing value. The index matrix is used to link the missing values in the data set with their corresponding rows in the replacement matrix.

- **call**: The matched function call.

- **model**: A list containing the names of the cluster variable, the target variables, and the predictor variables with fixed and random effects, respectively.

- **random.L1**: A character string denoting the handling of random residual covariance matrices (not used here; see `jomoImpute`).

- **prior**: The prior parameters used in the imputation model.

- **iter**: A list containing the number of burn-in iterations, the number of iterations between imputations, and the number of imputed data sets.

- **par.burnin**: A multi-dimensional array containing the parameters of the imputation model from the burn-in phase.

- **par.imputation**: A multi-dimensional array containing the parameters of the imputation model from the imputation phase.

### Note

For objects of class `mitml`, methods for the generic functions `print`, `summary`, and `plot` have been defined. `mitmlComplete` is used for extracting the imputed data sets.

### Author(s)

Simon Grund, Alexander Robitzsch, Oliver Luedtke
References

See Also

`jomoImpute, mitmlComplete, summary.mitml, plot.mitml`

Examples

# NOTE: The number of iterations in these examples is much lower than it
# should be! This is done in order to comply with CRAN policies, and more
# iterations are recommended for applications in practice!

data(studentratings)

# *** ........................................
# the 'type' interface
#
# * Example 1.1: 'ReadDis' and 'SES', predicted by 'ReadAchiev' and
# 'CognAbility', with random slope for 'ReadAchiev'

type <- c(-2,0,0,0,0,3,1,2,0)
names(type) <- colnames(studentratings)
type

imp <- panImpute(studentratings, type=type, n.burn=1000, n.iter=100, m=5)

# * Example 1.2: 'ReadDis' and 'SES' groupwise for 'FedState',
# and predicted by 'ReadAchiev'

type <- c(-2,-1,0,0,0,2,1,0,0)
names(type) <- colnames(studentratings)
type

imp <- panImpute(studentratings, type=type, n.burn=1000, n.iter=100, m=5)

# *** ........................................
# the 'formula' interface
#
# * Example 2.1: imputation of 'ReadDis', predicted by 'ReadAchiev'
# (random intercept)

fml <- ReadDis ~ ReadAchiev + (1|ID)
imp <- panImpute(studentratings, formula=fml, n.burn=1000, n.iter=100, m=5)

# ... the intercept can be suppressed using '0' or '-1' (here for fixed intercept)

fml <- ReadDis ~ 0 + ReadAchiev + (1|ID)
imp <- panImpute(studentratings, formula=fml, n.burn=1000, n.iter=100, m=5)
# Example 2.2: imputation of 'ReadDis', predicted by 'ReadAchiev'
# (random slope)

```r
fml <- readDis ~ readAchiev + (1+readAchiev|ID)
imp <- panImpute(studentratings, formula=fml, n.burn=1000, n.iter=100, m=5)
```

# Example 2.3: imputation of 'ReadDis', predicted by 'ReadAchiev',
# groupwise for 'FedState'

```r
fml <- readDis ~ readAchiev + (1|ID)
imp <- panImpute(studentratings, formula=fml, group="FedState", n.burn=1000,
                 n.iter=100, m=5)
```

# Example 2.4: imputation of 'ReadDis', predicted by 'ReadAchiev'
# including the cluster mean of 'ReadAchiev' as an additional predictor

```r
fml <- readDis ~ readAchiev + I(clusterMeans(readAchiev,ID)) + (1|ID)
imp <- panImpute(studentratings, formula=fml, n.burn=1000, n.iter=100, m=5)
```

# ... using 'save.pred' to save the calculated cluster means in the data set

```r
fml <- readDis ~ readAchiev + I(clusterMeans(readAchiev,ID)) + (1|ID)
imp <- panImpute(studentratings, formula=fml, n.burn=1000, n.iter=100, m=5,
                 save.pred=TRUE)
```

```r
head(mitmlComplete(imp,1))
```

---

**plot.mitml**

Print diagnostic plots

**Description**

Generates diagnostic plots for assessing the convergence and autocorrelation behavior of pan's and jomo's MCMC algorithms.

**Usage**

```r
# S3 method for class 'mitml'
plot(x, print=c("beta","beta2","psi","sigma"), pos=NULL, group="all",
     trace=c("imputation","burnin","all"), thin=1, smooth=3, n.Rhat=3,
     export=c("none","png","pdf"), dev.args=list(), ...)
```

**Arguments**

- `x` An object of class mitml as produced by panImpute and jomoImpute.
- `print` A character vector containing one or several of "beta", "beta2", "psi" or "sigma" denoting which parameters should be plotted. Default is to plot all parameters.
pos
Either NULL or an integer vector denoting a specific entry in either "beta", "beta2", "psi" or "sigma". Default is to NULL plotting all entries.

group
Either "all" or an integer denoting for which group plots should be generated. Used only when group-wise imputation was used. Default is to "all".

trace
One of "imputation", "burnin" or "all" denoting which part of the parameter chain should be used for the trace plot. Default is to "imputation", which plots only the iterations after burn-in.

thin
An integer denoting the thinning factor that is applied before plotting. Default is to 1, plotting the full chain.

smooth
A numeric value denoting the smoothing factor for the trend line in trace plots. Higher values correspond to less smoothing. Default is 3. If set to 0 or NULL, the trend line is suppressed.

n.Rhat
An integer denoting the number of sequences used for calculating the potential scale reduction factor. Default is 3.

export
(optional) A character string specifying if plots should be exported to file. If "png" or "pdf", then plots are printed into a folder named "mitmlPlots" in the current directory using either the png or pdf device. Default is to "none", which does not export files.

dev.args
(optional) A named list containing additional arguments that are passed to the graphics device.

... Parameters passed to the plotting functions.

Details
The `plot` method generates a series of plots for the parameters of the imputation model which can be used for diagnostic purposes. In addition, a short summary of the parameter chain is displayed.

Setting `print` to "beta", "beta2", "psi" and "sigma" will plot the fixed effects, the variances and covariances of random effects, and the variances and covariances of residuals, respectively. Here, "beta2" refers to the fixed effects for target variables at level 2 and is only used when imputations were carried out using a two-level model (`jomoImpute`). Each plotting window contains a trace plot (upper left), an autocorrelation plot (lower left), a kernel density approximation of the posterior distribution (upper right), and a posterior summary (lower right). The summary includes the following quantities:

**EAP:** Expected value a posteriori (i.e., the mean of the parameter chain)

**MAP:** Mode a posteriori (i.e., the mode of the parameter chain)

**SD:** Standard deviation of the parameter chain

**2.5%:** The 2.5% quantile of parameter values

**97.5%:** The 97.5% quantile of parameter values

**Rhat:** Estimated potential scale reduction factor ($\hat{R}$)

**ACF-k:** Smoothed autocorrelation at lag $k$, where $k$ is the number of iterations between imputations (see `summary.mitml`)
The `trace` and `smooth` arguments can be used to influence how the trace plot is drawn and what part of the chain should be used for it. The `thin` argument can be used for thinning the chain before plotting, in which case the number of data points is reduced in the trace plot, and the autocorrelation is calculated up to lag \( k/thin \) (see above). The `n.Rhat` argument controls the number of sequences that are used for calculating the potential scale reduction factor (\( \hat{R} \)) in each plot (see `summary.mitml`). Further arguments to the graphics device are supplied using the `dev.args` argument.

The `plot` function calculates and displays diagnostic information primarily for the imputation phase (i.e., for iterations after burn-in). This is the default in the `plot` function and the recommended method for most users. However, note that, when overriding the default using `trace="burnin"`, the posterior summary and the trace plots do not convey the necessary information to establish convergence. When `trace="all"`, the full chain is displayed with emphasis on the imputation phase, and the posterior summary is calculated based only on iterations after burn-in as recommended.

**Value**

None (invisible `NULL`).

**Note**

The plots are presented on-screen one at a time. To proceed with the next plot, the user may left-click in the plotting window or press the "enter" key while in the R console, depending on the operating system. No plots are displayed when exporting to file.

**Author(s)**

Simon Grund

**See Also**

`panImpute`, `jomoImpute`, `summary.mitml`

**Examples**

```r
## Not run:
data(studentratings)

# * Example 1: simple imputation

fml <- ReadDis + SES ~ ReadAchiev + (1|ID)
imp <- panImpute(studentratings, formula=fml, n.burn=1000, n.iter=100, m=5)

# plot all parameters (default)
plot(imp)

# plot fixed effects only
plot(imp, print="beta")

# export plots to file (using pdf device)
plot(imp, export="pdf", dev.args=list(width=9, height=4, pointsize=12))
```
# Example 2: groupwise imputation

```r
fml <- ReadDis + SES ~ ReadAchiev + (1|ID)
imp <- panImpute(studentratings, formula=fml, group=FedState, n.burn=1000, n.iter=100, m=5)

# plot fixed effects for all groups (default for 'group')
plot(imp, print="beta", group="all")

# plot fixed effects for first group only
plot(imp, print="beta", group=1)

## End(Not run)
```

---

**read.mitml**  
*Read mitml objects from file*

**Description**

This function loads *mitml* class objects from R binary formats (similar to *load*), usually produced by `write.mitml`.

**Usage**

```r
read.mitml(filename)
```

**Arguments**

`filename`  
Name of the file to read, to be specified with file extension (e.g., `.R`, `.Rdata`).

**Value**

Returns the saved *mitml* class object.

**Author(s)**

Simon Grund

**See Also**

`panImpute`, `jomoImpute`, `write.mitml`
sort.mitml.list

Examples

data(studentratings)

fml <- ReadDis + SES ~ ReadAchiev + (1|ID)
imp <- panImpute(studentratings, formula=fml, n.burn=1000, n.iter=100, m=5)

# write 'mitml' object
write.mitml(imp, filename="imputation.R")

# read previously saved 'mitml' object
previous.imp <- read.mitml("imputation.R")
class(previous.imp)
previous.imp

sort.mitml.list Sort a list of imputed data sets

Description

The functions sorts a list of multiply imputed data sets according to an R expression.

Usage

## S3 method for class 'mitml.list'
sort(x, decreasing=FALSE, by, ...)

Arguments

x A list of imputed data sets with class mitml.list as produced by mitmlComplete
(or similar).
decreasing Logical flag indicating if data sets should be sorted in decreasing (i.e., reversed)
order. Default is 'FALSE'.
by An R expression or a list of multiple expressions by which to sort the imputed
data sets (see Examples).
... Further arguments to 'order' (see Details).

Details

This function sorts a list of imputed data sets according to the R expression given in the by argument.
The function is similar to the order function for regular data sets and uses it internally. Note that
sorting is performed individually for each data set. Thus, the order of cases may differ across data
sets if the variables used for sorting contain different values.
Value

A list of imputed data sets with an additional class attribute `mitml.list`.

Author(s)

Simon Grund

Examples

data(studentratings)

```r
fml <- ReadDis + SES - ReadAchiev + (1|ID)
imp <- panImpute(studentratings, formula=fml, n.burn=1000, n.iter=100, m=5)

implist <- mitmlComplete(imp,"all")
```

`# * Example 1: sort by ID`

```r
sort(implist, by=ID)
```

`# * Example 2: sort by combination of variables`

```r
sort(implist, by=list(FedState,ID,-SES))
```

---

`studentratings`  
*Example data set on student’s ratings and achievement*

Description

This data set contains simulated data for students nested within schools, featuring students’ ratings of their teachers’ behavior (i.e., disciplinary problems in mathematics and reading class) and their general learning environment (school climate), as well as mathematics and reading achievement scores, and scores for socio-economic status and cognitive ability.

In addition, the data set features the ID of 50 different schools (i.e., clusters), the biological sex of all students, and a broad, additional grouping factor. Different amounts of missing data have been inserted into the data set in a completely random fashion.

Usage

```r
data(studentratings)
```

Format

A data frame containing 750 observations on 10 variables.
Subset a list of imputed data sets

Description
The functions can be used for creating subsets for a list of multiply imputed data sets.

Usage

```r
## S3 method for class 'mitml.list'
subset(x, subset, select, ...)
```

Arguments

- `x`: A list of imputed data sets with class `mitml.list` as produced by `mitmlComplete` (or similar).
- `subset`: An R expression by which to subset each data set.
- `select`: An R expression by which to select columns.
- `...`: Not being used.

Details
This function can be used to create subsets and select variables for a list of multiply imputed data sets according to the R expressions given in the `subset` and `select` arguments. The function is similar to and adapted from the `subset` function for regular data sets. Note that subsetting is performed individually for each data set. Thus, the cases included may differ across data sets if the variables used for subsetting contain different values.

Value
A list of imputed data sets with an additional class attribute `mitml.list`.

Author(s)
Simon Grund

Examples

```r
data(studentratings)
fml <- ReadDis + SES ~ ReadAchiev + (1|ID)
imp <- panImpute(studentratings, formula=fml, n.burn=1000, n.iter=100, m=5)
implist <- mitmlComplete(imp,"all")
```
# Example 1: subset by SES, select variables by name
subset(implist, SES < 25, select = c(ID, FedState, Sex, SES, ReadAchiev, ReadDis))

# Example 2: subset by FedState, select variables by column number
subset(implist, FedState == "SH", select = c(6:7, 9:10))

## Not run:
# Example 3: subset by ID and Sex
subset(implist, ID)

# Example 4: select variables by name range
subset(implist, select = ID:Sex)

## End(Not run)

summary.mitml  

**Summary measures for imputation models**

**Description**

Provides summary statistics and additional information on imputations in objects of class `mitml`.

**Usage**

```r
## S3 method for class 'mitml'
summary(object, n.Rhat=3, goodness.of.appr=FALSE, autocorrelation=FALSE, ...)
```

**Arguments**

- **object**: An object of class `mitml` as produced by `panImpute` and `jomoImpute`
- **n.Rhat**: (optional) An integer denoting the number of sequences used for calculating the potential scale reduction factor. Default is to 3.
- **goodness.of.appr**: (optional) A logical flag indicating if the goodness of approximation should be printed. Default is to FALSE (see details).
- **autocorrelation**: (optional) A logical flag indicating if the autocorrelation should be printed. Default is to FALSE (see details).
- **...**: Not being used.

**Details**

The `summary` method calculates summary statistics for objects of class `mitml` as produced by `panImpute` and `jomoImpute`. The output includes the potential scale reduction factor (PSRF, or $\hat{R}$) and (optionally) the goodness of approximation and autocorrelation.
The PSRF is calculated for each parameter of the imputation model and may be interpreted as a measure of convergence (Gelman and Rubin, 1992). Calculation of the PSRFs can be suppressed by setting \( n.Rhat=NULL \). The PSRFs are not computed from different chains, but by dividing each chain from the imputation phase into a number of sequences as denoted by \( n.Rhat \). This is slightly different from the original method proposed by Gelman and Rubin.

The goodness of approximation indicates what proportion of the posterior standard deviation is due to simulation error. For multiple imputation, the goodness of approximation is not essential; it should be considered only if posterior summaries, such as the EAP, are of interest.

The autocorrelation includes estimates of the autocorrelation in the parameter chains at lag 1 (i.e., for consecutive draws) and for lags \( k \) and \( 2k \), where \( k \) is the number of iterations between imputations. For lag \( k \) and \( 2k \), the autocorrelation is slightly smoothed to reduce the influence of noise on the estimates (see \texttt{plot.mitml}.

**Value**

Returns an object of class \texttt{summary.mitml}. A print method is used for better readable console output.

**Author(s)**

Simon Grund

**References**


**See Also**

\texttt{panImpute, jomoImpute, plot.mitml}

**Examples**

data(studentratings)

fml <- ReadDis + SES ~ ReadAchiev + (1|ID)
imp <- panImpute(studentratings, formula=fml, n.burn=1000, n.iter=100, m=5)

# print summary
summary(imp)
**Description**

Performs hypothesis tests for arbitrary functions of a parameter vector using the Delta method.

**Usage**

```r
testConstraints(model, qhat, uhat, constraints, method=c("D1","D2"), df.com=NULL)
```

**Arguments**

- **model**: A list of fitted statistical models (see examples).
- **qhat, uhat**: Two matrices/arrays or lists containing estimates of the parameter vector and its covariance matrix, respectively, for each imputed data set (see examples).
- **constraints**: A character vector specifying constraints or functions of the vector of model parameters to be tested.
- **method**: A character string denoting the method by which the test is performed. Can be either "D1" or "D2" (see details). Default is to "D1".
- **df.com** (optional): A single number or a numeric vector denoting the complete-data degrees of freedom for the hypothesis test. Only used if method= "D1".

**Details**

This function is similar in functionality to `testModels` but extended to arbitrary functions (or constraints) of the model parameters. The function is based on the Delta method (e.g., Casella & Berger, 2002) according to which any function of the parameters can be tested using Wald-like methods, assuming that their sampling distribution is approximately normal. It is assumed that the parameters can be extracted using `coef` and `vcov` methods from the fitted models (or similar; e.g., regression coefficients, fixed effects in multilevel models). In cases where this is not possible, hypothesis tests can be carried out using user-supplied matrices/arrays or lists (qhat and uhat, see examples).

Constraints and functions of the model parameters can be specified in the `constraints` argument. The constraints must be supplied as a character vector, where each string denotes a function or a constraint to be tested (see examples).

The Wald-like tests that are carried out by `testConstraints` can be aggregated across data sets with the two methods $D_1$ (Li, Raghunathan & Rubin, 1991) and $D_2$ (Li, Meng, Raghunathan & Rubin, 1991), where $D_1$ operates on the constrained estimates and standard errors, and $D_2$ operates on the Wald-statistics (for an explanation, see `testModels`). The pooled estimates and standard errors reported in the output are always based on $D_1$.

For $D_1$, the complete-data degrees of freedom can be adjusted for smaller samples by specifying `df.com`. 
Currently, the procedure supports statistical models that define `coef` and `vcov` methods (e.g., `lm`), multilevel models estimated with `lme4` or `nlme`, and GEEs estimated with `geepack`. The arguments `qhat` and `uhat` allow for more general hypothesis tests regardless of model class. Support for further models may be added in future releases.

**Value**

Returns a list containing the results of the model comparison, the constrained estimates and standard errors, and the relative increase in variance due to nonresponse (Rubin, 1987). A `print` method is used for better readable console output.

**Author(s)**

Simon Grund

**References**


**See Also**

`testModels, withNmitmlNlist`

**Examples**

data(studentratings)

```r
fml <- MathDis + ReadDis + SchClimate ~ (1|ID)
ip <- panImpute(studentratings, formula=fml, n.burn=1000, n.iter=100, m=5)

implist <- mitmlComplete(ip, print=1:5)

# fit simple regression model
fit.lm <- with(implist, lm(SchClimate ~ ReadDis + MathDis))

# apply Rubin's rules
testEstimates(fit.lm)

# Example 1: test 'identity' function of two parameters
# multi-parameter hypothesis test, equivalent to model comparison

cons <- c("ReadDis","MathDis")
testConstraints(fit.lm, constraints=cons)

# ... adjusting for finite samples
testConstraints(fit.lm, constraints=cons, df.com=749)
```
testEstimates: Compute final estimates and inferences

Description

Computes final parameter estimates and inferences from multiply imputed data sets.

Usage

testEstimates(model, qhat, uhat, var.comp=FALSE, df.com=NULL)

Arguments

- **model**: A list of fitted statistical models (see examples).
- **qhat, uhat**: Two matrices or lists containing point and variances estimates, respectively, for each imputed data set (see examples).
- **var.comp**: A logical flag indicating if estimates for variance components should be calculated. Default is to FALSE.
- **df.com**: (optional) A numeric vector denoting the complete-data degrees of freedom for the hypothesis test.
Details

This function calculates final parameter estimates and inferences as suggested by Rubin (1987, "Rubin’s rules") for each parameter of the fitted model. In other words, testEstimates aggregates
estimates and standard errors across multiply imputed data sets. The parameters and standard errors
can either be supplied as fitted statistical models (model), or as two matrices or lists (qhat and uhat,
see examples).

Rubin’s original method assumes that the complete-data degrees of freedom are infinite, which is
reasonable in larger samples. Alternatively, the degrees of freedom can be adjusted for smaller
samples by specifying df.com (Barnard & Rubin, 1999). The df.com argument can either be a
single number if the degrees of freedom are equal for all tests, or a numeric vector with one element
per test.

Using the var.comp argument, final estimates for variance components and related parameters can
be requested. These will be shown as a separate table within the console output. Accessing variance
components is highly dependent on the model being estimated and not implemented for all models.

Users may prefer calculating these estimates manually using with.mitml.list (see Example 3). No
inferences are calculated for variance components.

Currently, the procedure supports statistical models that define coef and vcov methods (e.g., lm),
multilevel models estimated with lme4 or nlme, and GEEs estimated with gee pack. The arguments
qhat and uhat allow for more general aggregation of parameter estimates regardless of model class.
Support for further models may be added in future releases.

Value

Returns a list containing the final parameter and inferences, the relative increase in variance due to
nonresponse, and the fraction of missing information (Rubin, 1987). A print method is used for
better readable console output.

Author(s)

Simon Grund

References

Biometrika, 86, 948-955.


See Also

with.mitml.list, confint.mitml.testEstimates

Examples

data(studentsratings)

fml <- ReadDis + SES ~ ReadAchiev + (1|ID)
imp <- panImpute(studentsratings, formula=fml, n.burn=1000, n.iter=100, m=5)
testModels <- mitmlComplete(imp, print=1:5)

# fit multilevel model using lme4
require(lme4)
fit.lmer <- with(implist, lmer(SES ~ (1|ID)))

# * Example 1: combine estimates using model recognition
# final estimates and inferences separately for each parameter (Rubin's rules)
testEstimates(fit.lmer)

# ... adjusted df for finite samples
testEstimates(fit.lmer, df.com=49)

# ... with additional table for variance components and ICCs
testEstimates(fit.lmer, var.comp=TRUE)

# * Example 2: combine estimates using matrices or lists
fit.lmer <- with(implist, lmer(SES ~ ReadAchiev + (1|ID)))
qhat <- sapply(fit.lmer, fixef)
what <- sapply(fit.lmer, function(x) diag(vcov(x)))
testEstimates(qhat=qhat, what=what)

testModels

Test multiple parameters and compare nested models

Description
Performs multi-parameter hypothesis tests for a vector of statistical parameters and compares nested statistical models obtained from multiply imputed data sets.

Usage

testModels(model, null.model, method=c("D1","D2","D3"),
use=c("wald","likelihood"), df.com=NULL)

Arguments

model A list of fitted statistical models (see examples).
null.model A list of fitted (more restrictive) statistical models.
method A character string denoting the method by which the test is performed. Can be either "D1", "D2" or "D3" (see details). Default is "D1".
use A character string denoting Wald- or likelihood-based based tests. Can be either "wald" or "likelihood". Only used if method="D2".
df.com (optional) A single number or a numeric vector denoting the complete-data degrees of freedom for the hypothesis test. Only used if method="D1".
testModels

Details

This function compares two nested statistical models which differ by one or more parameters. In other words, the function performs Wald-like and likelihood-ratio hypothesis tests for the statistical parameters by which the two models differ.

The general approach to Wald-like inference for multi-dimensional estimands was introduced Rubin (1987) and further developed by Li, Raghunathan and Rubin (1991). This procedure is commonly referred to as \( D_1 \) and can be used by setting `method = "D1"`. \( D_1 \) is the multi-parameter equivalent of `testEstimates`, that is, it tests multiple parameters simultaneously. For \( D_1 \), the complete-data degrees of freedom are assumed to be infinite, but they can be adjusted for smaller samples by supplying `dfcom` (Reiter, 2007).

An alternative method for Wald-like hypothesis tests was suggested by Li, Meng, Raghunathan and Rubin (1991). The procedure is often called \( D_2 \) and can be used by setting `method = "D2"`. \( D_2 \) calculates the Wald-test directly for each data set and then aggregates the resulting \( \chi^2 \) values. The source of these values is specified by the `use` argument. If `use = "wald"` (the default), then a Wald-like hypothesis test similar to \( D_1 \) is performed. If `use = "likelihood"`, then the two models are compared through their likelihood.

A third method relying on likelihood-based comparisons was suggested by Meng and Rubin (1992). This procedure is referred to as \( D_3 \) and can be used by setting `method = "D3"`. \( D_3 \) compares the two models by aggregating the likelihood-ratio test across multiply imputed data sets.

In general, Wald-like hypothesis tests (\( D_1 \) and \( D_2 \)) are appropriate if the parameters can be assumed to follow a multivariate normal distribution (e.g., regression coefficients, fixed effects in multilevel models). Likelihood-based comparisons (\( D_2 \) and \( D_3 \)) are also appropriate in such cases and may also be used for variance components.

The function supports different classes of statistical models depending on which method is chosen. \( D_1 \) supports quite general models as long as they define `coef` and `vcov` methods (or similar) for extracting the parameter estimates and their estimated covariance matrix. \( D_2 \) can be used for the same models (if `use = "wald"`, or alternatively, for models that define a `logLik` method (if `use = "likelihood"`). Finally, \( D_3 \) supports linear models and linear mixed-effects models with a single cluster variable as estimated by `lme4` or `nlme` (see Laird, Lange, & Stram, 1987). Support for other statistical models may be added in future releases.

Value

Returns a list containing the results of the model comparison, and the relative increase in variance due to nonresponse (Rubin, 1987). A `print` method is used for better readable console output.

Note

The \( D_3 \) method and the likelihood-based \( D_2 \) assume that models were fit using maximum likelihood (ML). Models fit using REML are automatically refit using ML.

Author(s)

Simon Grund
References


See Also

*anova.mitml.result, testEstimates, testConstraints, with.mitml.list*

Examples

data(studentratings)

fml <- ReadDis + SES ~ ReadAchiev + (1|ID)
imp <- panImpute(studentratings, formula=fml, n.burn=1000, n.iter=100, m=5)

implist <- mitmlComplete(imp, print=1:5)

# * Example 1: multiparameter hypothesis test for 'ReadDis' and 'SES'
# This tests the hypothesis that both effects are zero.
require(lme4)
fit0 <- with(implist, lmer(ReadAchiev ~ (1|ID), REML=FALSE))
fit1 <- with(implist, lmer(ReadAchiev ~ ReadDis + (1|ID), REML=FALSE))

# apply Rubin's rules
testEstimates(fit1)

# multiparameter hypothesis test using D1 (default)
testModels(fit1, fit0)

# ... adjusting for finite samples
testModels(fit1, fit0, df.com=47)

# ... using D2 ("wald", using estimates and covariance-matrix)
testModels(fit1, fit0, method="D2")

# ... using D2 ("likelihood", using likelihood-ratio test)
testModels(fit1, fit0, method="D2", use="likelihood")
# ... using D3 (likelihood-ratio test, requires ML fit)
testModels(fit1, fit0, method="D3")

## Not run:
## * Example 2: multiparameter test using D3 with nlme

## for D3 to be calculable, the 'data' argument for 'lme' must be
## can be constructed manually

require(nlme)
fit0 <- withC
lme(ReadAchiev~1, random=~1|ID,
data=data.frame(ReadAchiev,ID, method="ML"))
fit1 <- withC
lme(ReadAchiev ~ 1 + ReadDis, random=~ 1|ID,
data=data.frame(ReadAchiev,ReadDis,ID), method="ML")

# multiparameter hypothesis test using D3
testModels(fit1, fit0, method="D3")

## End(Not run)

---

### with.mitml.list

**Evaluate an expression in a list of imputed data sets**

---

**Description**

The functions `with` and `within` evaluate R expressions in a list of multiply imputed data sets.

**Usage**

```r
## S3 method for class 'mitml.list'
with(data, expr, ...)
## S3 method for class 'mitml.list'
within(data, expr, ignore=NULL, ...)
```

**Arguments**

- `data` A list of imputed data sets with class `mitml.list` as produced by `mitmlComplete`.
- `expr` An R expression to be evaluated for each data set.
- `ignore` A character vector denoting objects not to be saved.
- `...` Not being used.

**Details**

The two functions are defined as `with` and `within` methods for objects of class `mitml.list`. Both `with` and `within` evaluate an R expression in each of the imputed data sets. However, the two functions return different values: `with` returns the evaluated expression, whereas `within` returns the resulting data sets. The `ignore` argument may be used to declare objects that are not to be saved within `within`.
Value

with: Returns the evaluated expression from each data set as a list (class mitml.result). This is useful for fitting statistical models to multiply imputed data. The list of fitted models can be analyzed using testEstimates, testModels, testConstraints, or anova.

within: Evaluates the R expression for each data set and returns the altered data sets as a list mitml.list. This is useful for manipulating the data prior to data analysis (e.g., centering, calculating cluster means, etc.).

Author(s)

Simon Grund

See Also

mitmlComplete, anova.mitml.result, testEstimates, testModels, testConstraints

Examples

data(studentratings)

fm1 <- ReadDis + SES ~ ReadAchiev + (1|ID)
imp <- panImpute(studentratings, formula=fm1, n.burn=1000, n.iter=100, m=5)

implist <- mitmlComplete(imp, print=1:5)

# * Example 1: data manipulation

# calculate and save cluster means
new1.implist <- within(implist, Means.ReadAchiev <- clusterMeans(ReadAchiev, ID))

# center variables, calculate interaction terms, ignore byproducts
new2.implist <- within(implist,{
    M.SES <- mean(SES)
    M.CognAbility <- mean(CognAbility)
    C.SES <- SES - M.SES
    C.CognAbility <- CognAbility - M.CognAbility
    SES.CognAbility <- C.SES * C.CognAbility
    }, ignore=c("M.SES", "M.CognAbility"))

# * Example 2: fitting statistical models

# fit regression model
fit.lm <- with(implist, lm(ReadAchiev ~ ReadDis))

# fit multilevel model using lme
require(lme4)
fit.lmer <- with(implist, lmer(ReadAchiev ~ ReadDis + (1|ID)))

# * Example 3: manual extraction of variance estimates
require(lme4)
fit.lmer <- with(implist, lmer(SES ~ (1|ID)))
write.mitml

# extract level-1 and level-2 variances
var.l1 <- sapply(fit.lmer, function(z) attr(VarCorr(z),"sc")^2)
var.l2 <- sapply(fit.lmer, function(z) VarCorr(z)$ID[1,1])

# calculate final estimate of the intraclass correlation
ICC <- mean( var.l2 / (var.l2+var.l1) )

---

write.mitml Write mitml objects to file

Description

This function saves objects of class mitml in R binary formats (similar to ?save).

Usage

write.mitml(x, filename, drop=FALSE)

Arguments

x An object of class mitml as produced by panImpute and jomoImpute.
filename Name of the destination file, to be specified with file extension (e.g., .R, .Rdata).
drop Logical flag indicating if the parameters of the imputation model should be dropped in favor for lower file size. Default is to FALSE.

Value

None (invisible NULL).

Author(s)

Simon Grund

See Also

panImpute, jomoImpute, read.mitml

Examples

data(studentratings)

fml <- ReadDis + SES ~ ReadAchiev + (1|ID)
imp <- panImpute(studentratings, formula=fml, n.burn=1000, n.iter=100, m=5)

# write full 'mitml' object (default)
write.mitmlMplus(imp, filename="imputation.R")

# drop parameters of the imputation model
write.mitml(imp, filename="imputation.R", drop=TRUE)

write.mitmlMplus Write mitml objects to Mplus format

Description

Saves objects of class mitml as a series of text files which can be processed by the statistical software Mplus (Muthen & Muthen, 2012).

Usage

write.mitmlMplus(x, filename, suffix="list", sep="\t", dec=".", na.value=-999)

Arguments

x An object of class mitml or mitml.list as produced by panImpute, jomoImpute, mitmlComplete, or similar).
filename Basic file name for the text files containing the imputed data sets, to be specified without file extension.
suffix File name suffix for the index file.
sep The field separator.
dec The decimal separator.
na.value A numeric value coding the missing data in the resulting data files.

Details

The Mplus format for multiply imputed data sets comprises a series of text files, each containing one imputed data set, and an index file containing the names of all data files. During export, factors and character variables are converted to numeric. Therefore, write.mitmlMplus produces a log file which contains information about the data set and the factors that have been converted.

In addition, a basic Mplus input file is generated that can be used for setting up subsequent analysis models.

Value

None (invisible NULL).

Author(s)

Simon Grund
References


See Also

`panImpute`, `jomoImpute`, `mitmlComplete`

Examples

```r
data(studentratings)

fml <- ReadDis + SES ~ ReadAchiev + (1|ID)
imp <- panImpute(studentratings, formula=fml, n.burn=1000, n.iter=100, m=5)

# write imputation files, index file, and log file
write.mitmlplus(imp, filename="imputation", suffix="list", na.value=-999)
```

**Description**

`write.mitmlSAV` Write mitml objects to native SPSS format

Saves objects of class mitml in the .sav format used by the statistical software SPSS (IBM Corp., 2013). The function serves as a front-end for `write_sav` from the haven package.

**Usage**

`write.mitmlSAV(x, filename)`

**Arguments**

- `x` An object of class mitml or mitml.list as produced by `panImpute`, `jomoImpute`, `mitmlComplete`, or similar.
- `filename` Name of the destination file, to be specified with or without file extension. The file extension (.sav) is appended if necessary.

**Details**

This function exports multiply imputed data sets to a single .sav file, in which an Imputation_variable separates the original data and the various imputed data sets. Thus, `write.mitmlSAV` exports directly to the native SPSS format.

Alternatively, `write.mitmlSPSS` may be used for creating separate text and SPSS syntax files; an option that offers more control over the data format.
Value

None (invisible NULL).

Author(s)

Simon Grund

References


See Also

panImpute, jomoImpute, mitmlComplete, write.mitmlSPSS

Examples

data(studentratings)

fml <- ReadDis + SES ~ ReadAchiev + (1|ID)
imp <- panImpute(studentratings, formula=fml, n.burn=1000, n.iter=100, m=5)

# write data file and SPSS syntax
write.mitmlSAV(imp, filename="imputation")

write.mitmlSPSS Write mitml objects to SPSS compatible format

Description

Saves objects of class mitml as a text and a syntax file which can be processed by the statistical software SPSS (IBM Corp., 2013).

Usage

write.mitmlSPSS(x, filename, sep="\t", dec=".", na.value=-999, syntax=TRUE, locale=NULL)

Arguments

x An object of class mitml or mitml.list as produced by panImpute, jomoImpute, mitmlComplete, or similar.
filename Basic file name of the data and syntax files, to be specified without file extension.
sep The field separator.
dec The decimal separator.
write.mitmlSPSS

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<th>Description</th>
</tr>
</thead>
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<tr>
<td>na.value</td>
<td>A numeric value coding the missing data in the resulting data file.</td>
</tr>
<tr>
<td>syntax</td>
<td>A logical flag indicating if an SPSS syntax file should be generated. This file contains instructions for SPSS for reading in the data file. Default is to TRUE.</td>
</tr>
<tr>
<td>locale</td>
<td>(optional) A character string specifying the localization to be used in SPSS (e.g., &quot;en_US&quot;, &quot;de_DE&quot;). This argument may be specified if SPSS reads the data incorrectly due to conflicting locale settings.</td>
</tr>
</tbody>
</table>

Details

Multiply imputed data sets in SPSS are contained in a single file, in which an Imputation_variable separates the original data and the various imputed data sets. During export, factors are converted to numeric, whereas character variables are left "as is".

By default, write.mitmlSPSS generates a raw text file containing the data, along with a syntax file containing instructions for SPSS. This syntax file mimics SPSS’s functionality to read text files but circumvents certain problems that may occur when using the GUI. In order to read in the data, the syntax file must be opened and executed using SPSS. The syntax file may be altered manually if problems occur, for example, if the file path of the data file is not correctly represented in the syntax.

Alternatively, write.mitmlSAV may be used for exporting directly to the SPSS native .sav format. However, this may offer less control over the data format.

Value

None (invisible NULL).

Author(s)

Simon Grund

References


See Also

panImpute, jomoImpute, mitmlComplete, write.mitmlSAV

Examples

data(studentratings)

fml <- ReadDis + SES ~ ReadAchiev + (1|ID)
imp <- panImpute(studentratings, formula=fml, n.burn=1000, n.iter=100, m=5)

# write data file and SPSS syntax
write.mitmlSPSS(imp, filename="imputation", sep="\t", dec=".", na.value=-999, locale="en_US")
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