

Package ‘simsem’

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Type Package

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Author Sunthud Pornprasertmanit <psunthud@ku.edu>, Patrick Miller
<patr1ckm@ku.edu>, Alexander Schoemann <schoemann@ku.edu>

Maintainer Sunthud Pornprasertmanit <psunthud@ku.edu>

Depends R(>= 2.15), methods, lavaan, MASS

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Description This package can be used to generate data using the structural equation modeling framework. This package is tailored to use those simulated data for various purposes, such as model fit evaluation, power analysis, or missing data handling and planning.

License GPL (>= 2)

LazyLoad yes

URL <http://www.simsem.org>

R topics documented:

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| | |
|---------|------------|
| analyze | <i>TBA</i> |
|---------|------------|

Description

TBA

Usage

```
analyze(model, data, package="lavaan", miss=NULL, indLab=NULL,
aux=NULL, ...)
```

Arguments

| | |
|---------|-----|
| model | TBA |
| data | TBA |
| package | TBA |
| miss | TBA |

| | |
|--------|-----|
| indLab | TBA |
| aux | TBA |
| ... | TBA |

Value

TBA

Author(s)

Patrick Miller (Univeristy of Notre Dame; <pmille13@nd.edu>) Sunthud Pornprasertmanit (University of Kansas; <psunthud@ku.edu>)

Examples

```
loading <- matrix(0, 6, 2)
loading[1:3, 1] <- NA
loading[4:6, 2] <- NA
LY <- bind(loading, 0.7)

latent.cor <- matrix(NA, 2, 2)
diag(latent.cor) <- 1
RPS <- binds(latent.cor, 0.5)

RTE <- binds(diag(6))

VY <- bind(rep(NA,6),2)

CFA.Model <- model(LY = LY, RPS = RPS, RTE = RTE, modelType = "CFA")

dat <- generate(CFA.Model,200)
out <- analyze(CFA.Model,dat)
```

| | |
|-------|---|
| anova | <i>Provide a comparison of nested models and nonnested models across replications</i> |
|-------|---|

Description

This function will provide averages of model fit statistics and indices for nested models. It will also provide average differences of fit indices and power for likelihood ratio tests of nested models.

Arguments

| | |
|--------|--|
| object | SimResult object being described. Currently at least two objects must be included as arguments |
| ... | any additional arguments, such as additional objects or for the function with result object |

Value

A data frame that provides the statistics described above from all parameters. For using with `linkS4class{SimResult}`, the result is a list with two or three elements:

- `summary`: Average of fit indices across all replications
- `diff`: Average of the differences in fit indices across all replications
- `varyParam`: The statistical power of chi-square difference test given values of varying parameters (such as sample size or percent missing)

Author(s)

Alexander M. Schoemann (University of Kansas; <schoemann@ku.edu>), Sunthud Pornprasertmanit (University of Kansas; <psunthud@ku.edu>)

See Also

[SimResult](#) for the object input

Examples

```
loading1 <- matrix(0, 6, 1)
loading1[1:6, 1] <- NA
loading2 <- loading1
loading2[6,1] <- 0
LX1 <- bind(loading1, 0.7)
LX2 <- bind(loading2, 0.7)
RPH <- binds(diag(1))
RTD <- binds(diag(6))
CFA.Model1 <- model(LY = LX1, RPS = RPH, RTE = RTD, modelType="CFA")
CFA.Model2 <- model(LY = LX2, RPS = RPH, RTE = RTD, modelType="CFA")

# We make the examples running only 5 replications to save time.
# In reality, more replications are needed.
# Need to make sure that both simResult calls have the same seed!
Output1 <- sim(5, n=500, model=CFA.Model1, generate=CFA.Model1, seed=123567)
Output2 <- sim(5, n=500, model=CFA.Model2, generate=CFA.Model1, seed=123567)
anova(Output1, Output2)

Output1b <- sim(NULL, n=seq(50, 500, 50), model=CFA.Model1, generate=CFA.Model1, seed=123567)
Output2b <- sim(NULL, n=seq(50, 500, 50), model=CFA.Model2, generate=CFA.Model1, seed=123567)
anova(Output1b, Output2b)
```

bind

Specify matrices for Monte Carlo simulation of structural equation models

Description

Create [SimMatrix](#) or [SimVector](#) object that specifies

1. Pattern of fixed/freed parameters for analysis
2. Population parameter values for data generation

3. Any model misspecification (true population parameter is different than the one specified) for these parameters.

Each matrix in the Lisrel-style notation is specified in this way (e.g. LY, PS, and TE) and is used to create a model analysis template and a data generation template for simulation through the `model` function.

Usage

```
bind(free = NULL, popParam = NULL, misspec = NULL, symmetric = FALSE)
binds(free = NULL, popParam = NULL, misspec = NULL, symmetric = TRUE)
```

Arguments

| | |
|-----------|---|
| free | Required matrix or vector where each element represents a fixed or freed parameter used for analysis with structural equation models. Parameters can be freed by setting the corresponding element in the matrix to NA, and can be fixed by setting the value of the element to any number (e.g. 0). Parameters can be labeled using any character string. Any labeled parameter is considered to be free, and parameters with identical labels will be constrained to equality for analysis. |
| popParam | Optional matrix or vector of identical dimension to the free matrix whose elements contain population parameter values for data generation in simulation. For simulation, each free parameter requires a population parameter value, which is a quoted numeric value. Parameters that don't have population values are left as empty strings. Population parameters can also be drawn from a distribution. This is done by wrapping a call to create 1 value from an existing random generation function in quotes: e.g. "runif(1,0,1)", "rnorm(1,0,.01)". Every replication in the simulation will draw a parameter value from this distribution. The function checks that what is quoted is valid R. If a random population parameter is constrained to equality in the free matrix, <i>each drawn population parameter value will be the same</i> . More details on data generation is available in <code>?generate</code> , <code>?createData</code> , and <code>?draw</code> . To simplify the most common case, <code>popParam</code> can take 1 value or distribution and create a matrix or vector that assigns that population parameter or distribution to all freed parameters. These population values are used as starting values for analysis by default. |
| misspec | Optional matrix or vector of identical dimension to the free matrix whose elements contain population parameter values for specifying misspecification. Elements of the misspec matrix contain population parameters that are added to parameters that are fixed or have an existing population value. These parameters are also quoted numeric strings, and can optionally be drawn from distributions as described above. To simplify the most common case, <code>misspec</code> can take 1 value or distribution and create a matrix or vector that assigns that value or distribution to all previously specified fixed parameters. Details about misspecification are included in the data generation functions. |
| symmetric | Set as TRUE if the matrix created is symmetric (RPS/PS, RTE/TE). The function <code>binds</code> can also be used, which defaults to <code>symmetric = TRUE</code> |

Details

Bind is the first step in the `bind -> model -> sim` workflow of *simsem*, and this document outlines the user interface or language used to describe these simulations. This interface, while complex,

enables a wide array of simulation specifications for structural equation models by building on LISREL-style parameter specifications.

In simulations supported by *simsem*, a given parameter may be either fixed or freed for analysis, but may optionally also have a population value or distribution for data generation, or a value or distribution of misspecification. The purpose of `bind` is to stack these multiple meanings of a parameter into an object recognized by *simsem*, a `SimMatrix`. Each matrix in the Lisrel notation (e.g. LY, PS, TE, BE) becomes a `SimMatrix`, and is passed to the function `model`, which builds the data generation template and an analysis template (a lavaan parameter table), collectively forming a `SimSem` object, which can be passed to the function `sim` for simulation.

Value

`SimMatrix` or `SimVector` object that used for model specification for analysis and data generation in *simsem*.

Author(s)

Patrick Miller (Univeristy of Notre Dame; <pmille13@nd.edu>), Sunthud Pornprasertmanit (University of Kansas; <psunthud@ku.edu>)

See Also

- `model` To combine `simMatrix` objects into a complete data analysis and data generation template, which is a `SimSem` object
- `generate` To generate data using the *simsem* template.
- `analyze` To analyze real or generated data using the *simsem* template.

Examples

```
loading <- matrix(0, 6, 2)
loading[1:3, 1] <- NA
loading[4:6, 2] <- NA
loadingValues <- matrix(0, 6, 2)
loadingValues[1:3, 1] <- 0.7
loadingValues[4:6, 2] <- 0.7
LY <- bind(loading, loadingValues)
summary(LY)

# Set both factor correlations to .05
latent.cor <- matrix(NA, 2, 2)
diag(latent.cor) <- 1
RPS <- binds(latent.cor, 0.5)

# Misspecify all error covarainces
error.cor <- matrix(0, 6, 6)
diag(error.cor) <- NA
RTE <- binds(error.cor, 1, "runif(1, -.05, .05)")
```

| | |
|----------|------------------------------------|
| bindDist | Create a data distribution object. |
|----------|------------------------------------|

Description

Create a data distribution object

Usage

```
bindDist(margins, ..., p = NULL, keepScale = TRUE, reverse = FALSE)
```

Arguments

| | |
|-----------|---|
| margins | A character vector specifying all the marginal distributions. See the description of margins attribute of the Mvdc function for further details. |
| ... | A list whose each component is a list of named components, giving the parameter values of the marginal distributions. See the description of paramMargins attribute of the Mvdc function for further details. |
| p | Number of variables. If only one distribution object is listed, the p will make the same distribution objects for all variables. |
| keepScale | A vector representing whether each variable is transformed its mean and standard deviation or not. If TRUE, transform back to retain the mean and standard deviation of a variable equal to the model implied mean and standard deviation (with sampling error) |
| reverse | A vector representing whether each variable is mirrored or not. If TRUE, reverse the distribution of a variable (e.g., from positive skewed to negative skewed. If one logical value is specified, it will apply to all variables. |

Value

[SimDataDist](#) that saves analysis result from simulate data.

Author(s)

Sunthud Pornprasertmanit (University of Kansas; <psunthud@ku.edu>)

See Also

- [SimResult](#) for the type of resulting object

Examples

```
d1 <- list(df=2)
d2 <- list(df=3)
d3 <- list(df=4)
d4 <- list(df=5)
d5 <- list(df=3)
d6 <- list(df=4)
d7 <- list(df=5)
d8 <- list(df=6)
```

```
dist <- bindDist(c(rep("t", 4), rep("chisq", 8)), d1, d2, d3, d4, d5, d6, d7, d8, d5, d6, d7, d8)
```

| | |
|-------|--|
| clean | <i>Extract only converged replications in the result objects</i> |
|-------|--|

Description

Extract only the replications that are convergent in all supplied result objects ([SimResult](#))

Usage

```
clean(...)
```

Arguments

... The target result objects ([SimResult](#))

Value

The cleaned result objects

Author(s)

Sunthud Pornprasertmanit (University of Kansas; <psunthud@ku.edu>)

Examples

```
# No example
```

| | |
|----------------|---|
| cleanSimResult | <i>Extract only converged replications in the result object</i> |
|----------------|---|

Description

Extract only the replications that are convergent in a result object ([SimResult](#))

Usage

```
cleanSimResult(object, converged=NULL)
```

Arguments

| | |
|-----------|---|
| object | The target result object (SimResult) |
| converged | The replications to be extracted. If NULL, the converged slot in the result object will be used |

Value

The cleaned result object

Author(s)

Sunthud Pornprasertmanit (University of Kansas; <psunthud@ku.edu>)

Examples

```
# No example
```

| | |
|-----------------|---|
| continuousPower | <i>Find power of model parameters when simulations have randomly varying parameters</i> |
|-----------------|---|

Description

A function to find the power of parameters in a model when one or more of the simulations parameters vary randomly across replications.

Usage

```
continuousPower(simResult, contN = TRUE, contMCAR = FALSE, contMAR = FALSE,
  contParam = NULL, alpha = .05, powerParam = NULL, pred = NULL)
```

Arguments

| | |
|------------|--|
| simResult | SimResult that includes at least one randomly varying parameter (e.g. sample size, percent missing, model parameters) |
| contN | Logical indicating if N varies over replications. |
| contMCAR | Logical indicating if the percentage of missing data that is MCAR varies over replications. |
| contMAR | Logical indicating if the percentage of missing data that is MAR varies over replications. |
| contParam | Vector of parameters names that vary over replications. |
| alpha | Alpha level to use for power analysis. |
| powerParam | Vector of parameters names that the user wishes to find power for. This can be a vector of names (e.g., "LY1_1", "LY2_2"), or the name of a matrix (e.g. "PS"), if the name of a matrix is used power for all parameters in that matrix will be returned. If parameters are not specified, power for all parameters in the model will be returned. |
| pred | A list of varying parameter values that users wish to find statistical power from. |

Details

A common use of simulations is to conduct power analyses, especially when using SEM (Muthen & Muthen, 2002). Here, researchers select values for each parameter and a sample size and run a simulation to determine power in those conditions (the proportion of generated datasets in which a particular parameter of interest is significantly different from zero). To evaluate power at multiple sample sizes, one simulation for each sample size must be run. By continuously varying sample size across replications, only a single simulation is needed. In this simulation, the sample size for each replication varies randomly across plausible sample sizes (e.g., sample sizes between 200 and 500). For each replication, the sample size and significance of each parameter (0 = not significant, 1 = significant) are recorded. When the simulation is complete, parameter significance is regressed on sample size using logistic regression. For a given sample size, the predicted probability from the logistic regression equation is the power to detect an effect at that sample size. This approach can be extended to other randomly varying simulation parameters such as the percentage of missing data, and model parameters.

Value

Data frame containing columns representing values of the randomly varying simulation parameters, and power for model parameters of interest.

Author(s)

Alexander M. Schoemann (University of Kansas; <schoemann@ku.edu>), Sunthud Pornprasertmanit (University of Kansas; <psunthud@ku.edu>)

References

Muthen, L. K., & Muthen, B. O. (2002). How to use a Monte Carlo study to decide on sample size and determine power. *Structural Equation Modeling*, 4, 599-620.

See Also

- [SimResult](#) to see how to create a simResult object with randomly varying parameters.

Examples

```
## Not run:
# Specify Sample Size by n
loading <- matrix(0, 6, 1)
loading[1:6, 1] <- NA
LX <- bind(loading, 0.7)
RPH <- binds(diag(1))
RTD <- binds(diag(6))
CFA.Model <- model(LY = LX, RPS = RPH, RTE = RTD, modelType="CFA")
dat <- generate(CFA.Model, 50)
out <- analyze(CFA.Model, dat)

# We will use only 5 replications to save time.
# In reality, more replications are needed.

# Specify both sample size and percent missing completely at random

Output <- sim(NULL, CFA.Model, n=seq(100, 200, 20), pmMCAR=c(0, 0.1, 0.2))
summary(Output)

Cpow <- continuousPower(Output, contN = TRUE, contMCAR = TRUE)
Cpow

Cpow2 <- continuousPower(Output, contN = TRUE, contMCAR = TRUE, pred=list(N = 200, pmMCAR = 0.3))
Cpow2

## End(Not run)
```

Description

This function can be used to create data from a set of parameters created from [draw](#), called a code-paramSet. This function is used internally to create data, and is available publicly for accessibility and debugging.

Usage

```
createData(paramSet, n, indDist=NULL, sequential=FALSE, facDist=NULL,
  errorDist=NULL, indLab=NULL, modelBoot=FALSE, realData=NULL)
```

Arguments

| | |
|------------|---|
| paramSet | Set of drawn parameters from draw . |
| n | Integer of desired sample size. |
| indDist | A SimDataDist object or list of objects for a distribution of indicators. If one object is passed, each indicator will have the same distribution. Use when sequential is FALSE. |
| sequential | If TRUE, use a sequential method to create data such that the data from factor are generated first and apply to a set of equations to obtain the data of indicators. If FALSE, create data directly from model-implied mean and covariance of indicators. |
| facDist | A SimDataDist object or list of objects for the distribution of factors. If one object is passed, all factors will have the same distribution. Use when sequential is TRUE. |
| errorDist | An object or list of objects of type SimDataDist indicating the distribution of errors. If a single SimDataDist is specified, each error will be generated with that distribution. |
| indLab | A vector of indicator labels. When not specified, the variable names are x1, x2, ... xN. |
| modelBoot | When specified, a model-based bootstrap is used for data generation. See details for further information. This argument requires real data to be passed to readData. |
| realData | A data.frame containing real data. The data generated will follow the distribution of this data set. |

Details

This function will use `mvrnorm` function in MASS package to create data from model implied covariance matrix if the data distribution object ([SimDataDist](#)) is not specified. If the data distribution object is specified, the Gaussian copula model is used. See [SimDataDist](#) for further details. For the model-based bootstrap, the transformation proposed by Yung & Bentler (1996) is used. This procedure is the expansion from the Bollen and Stine (1992) bootstrap including a mean structure. The model-implied mean vector and covariance matrix with trivial misspecification will be used in the model-based bootstrap if `misspec` is specified. See page 133 of Bollen and Stine (1992) for a reference.

Internally, parameters are first drawn, and data is then created from these parameters. Both of these steps are available via the [draw](#) and [createData](#) functions respectively.

Value

A data.frame containing simulated data from the data generation template. A variable "group" is appended indicating group membership.

Author(s)

Sunthud Pornprasertmanit (University of Kansas; <psunthud@ku.edu>), Patrick Miller (University of Notre Dame; <pmille13@nd.edu>)

References

- Bollen, K. A., & Stine, R. A. (1992). Bootstrapping goodness-of-fit measures in structural equation models. *Sociological Methods and Research*, 21, 205-229.
- Yung, Y.-F., & Bentler, P. M. (1996). Bootstrapping techniques in analysis of mean and covariance structures. In G. A. Marcoulides & R. E. Schumacker (Eds.), *Advanced structural equation modeling: Issues and techniques* (pp. 195-226). Mahwah, NJ: Erlbaum.

Examples

```
loading <- matrix(0, 6, 2)
loading[1:3, 1] <- NA
loading[4:6, 2] <- NA
LY <- bind(loading, 0.7)

latent.cor <- matrix(NA, 2, 2)
diag(latent.cor) <- 1
RPS <- binds(latent.cor, 0.5)

RTE <- binds(diag(6))

VY <- bind(rep(NA,6),2)

CFA.Model <- model(LY = LY, RPS = RPS, RTE = RTE, modelType = "CFA")

param <- draw(CFA.Model)

# Generate data from the first group in the paramList.
dat <- createData(param[[1]], n = 200)
```

draw

Draw parameters from a [SimSem](#) object.**Description**

This function draws parameters from a [SimSem](#) template, for debugging or other use. Used internally to create data. Data can be created in one step from a [SimSem](#) object using [generate](#).

Usage

```
draw(model, maxDraw=50, misfitBounds=NULL, averageNumMisspec=FALSE,
      optMisfit = NULL, optDraws=50, misfitType="f0", misfitOut=FALSE)
```

Arguments

| | |
|---------|---|
| model | A SimSem object. |
| maxDraw | Integer specifying the maximum number of attempts to draw a valid set of parameters (no negative error variance, standardized coefficients over 1). |

| | |
|--------------------------------|---|
| <code>misfitBounds</code> | Vector that contains upper and lower bounds of the misfit measure. Sets of parameters drawn that are not within these bounds are rejected. |
| <code>averageNumMisspec</code> | TRUE or FALSE. ?? |
| <code>optMisfit</code> | Character vector of either "min" or "max" indicating either maximum or minimum optimized misfit. If not null, the set of parameters out of the number of draws in "optDraws" that has either the maximum or minimum misfit of the given misfit type will be returned. |
| <code>optDraws</code> | Number of parameter sets to draw if <code>optMisfit</code> is not null. The set of parameters with the maximum or minimum misfit will be returned. |
| <code>misfitType</code> | Character vector indicating the fit measure used to assess the misfit of a set of parameters. Can be "f0", "rmsea", "srmr", or "all". |
| <code>misfitOut</code> | Argument specifying how the misfit is returned. |

Value

Nested list of drawn parameters in the form `[[Group]][[param,misspec,misOnly]][[SimMatrix]]`. E.g. The LY parameter matrix of the first group would be indexed as `obj[[1]]$param$LY`. The values in `$param` are the raw parameter values with no misspecification. The values in `$misspec` are raw parameter values + misspecification. The values in `$misOnly` are only the misspecification values.

Author(s)

Sunthud Pornprasertmanit (University of Kansas; <psunthud@ku.edu>), Patrick Miller (Univeristy of Notre Dame; <pmille13@nd.edu>)

See Also

[createData](#) To generate random data using a set of parameters from [draw](#)

Examples

```
loading <- matrix(0, 6, 2)
loading[1:3, 1] <- NA
loading[4:6, 2] <- NA
LY <- bind(loading, 0.7)

latent.cor <- matrix(NA, 2, 2)
diag(latent.cor) <- 1
RPS <- binds(latent.cor, 0.5)

RTE <- binds(diag(6))

VY <- bind(rep(NA,6),2)

CFA.Model <- model(LY = LY, RPS = RPS, RTE = RTE, modelType = "CFA")

param <- draw(CFA.Model)
```

estmodel

*Shortcut for data analysis template for simulation.***Description**

Creates a data analysis template (lavaan parameter table) for simulations with structural equation models based on Y-side LISREL design matrices. Each corresponds to a LISREL matrix, but must be a matrix or a vector. In addition to the usual Y-side matrices in LISREL, both PS and TE can be specified using correlations (RPS, RTE) and scaled by a vector of residual variances (VTE, VPS) or total variances (VY, VE). Multiple groups are supported by passing lists of matrices or vectors to arguments, or by specifying the number of groups.

Usage

```
estmodel(LY = NULL, PS = NULL, RPS = NULL, TE = NULL, RTE = NULL, BE = NULL,
VTE = NULL, VY = NULL, VPS = NULL, VE=NULL, TY = NULL, AL = NULL,
MY = NULL, ME = NULL, modelType, indLab=NULL, facLab=NULL, groupLab="group",
ngroups=1, smartStart=TRUE)
estmodel.cfa(LY = NULL, PS = NULL, RPS = NULL, TE = NULL, RTE = NULL, VTE = NULL,
VY = NULL, VPS = NULL, VE=NULL, TY = NULL, AL = NULL, MY = NULL, ME = NULL,
indLab=NULL, facLab=NULL, groupLab="group", ngroups=1, smartStart=TRUE)
estmodel.path(PS = NULL, RPS = NULL, BE = NULL, VPS = NULL, VE=NULL, AL = NULL,
ME = NULL, indLab=NULL, facLab=NULL, groupLab="group", ngroups=1, smartStart=TRUE)
estmodel.sem(LY = NULL, PS = NULL, RPS = NULL, TE = NULL, RTE = NULL, BE = NULL,
VTE = NULL, VY = NULL, VPS = NULL, VE=NULL, TY = NULL, AL = NULL, MY = NULL,
ME = NULL, indLab=NULL, facLab=NULL, groupLab="group", ngroups=1, smartStart=TRUE)
```

Arguments

| | |
|-----|---|
| LY | Factor loading matrix from endogenous factors to Y indicators (need to be a matrix or a list of matrices). |
| PS | Residual covariance matrix among endogenous factors (need to be a symmetric matrix or a list of symmetric matrices). |
| RPS | Residual correlation matrix among endogenous factors (need to be a symmetric matrix or a list of symmetric matrices). |
| TE | Measurement error covariance matrix among Y indicators (need to be a symmetric matrix or a list of symmetric matrices). |
| RTE | Measurement error correlation matrix among Y indicators (need to be a symmetric matrix or a list of symmetric matrices). |
| BE | Regression coefficient matrix among endogenous factors (need to be a matrix or a list of matrices). |
| VTE | Measurement error variance of indicators (need to be a vector or a list of vectors). |
| VY | Total variance of indicators (need to be a vector or a list of vectors). NOTE: Either measurement error variance or indicator variance is specified. Both cannot be simultaneously specified. |
| VPS | Residual variance of factors (need to be a vector or a list of vectors). |

| | |
|------------|---|
| VE | Total variance of factors (need to be a vector or a list of vectors). NOTE: Either residual variance of factors or total variance of factors is specified. Both cannot be simultaneously specified. |
| TY | Measurement intercepts of Y indicators. (need to be a vector or a list of vectors). |
| AL | Endogenous factor intercept (need to be a vector or a list of vectors). |
| MY | Overall Y indicator means. (need to be a vector or a list of vectors). NOTE: Either measurement intercept of indicator mean can be specified. Both cannot be specified simultaneously. |
| ME | Total mean of endogenous factors (need to be a vector or a list of vectors). NOTE: Either endogenous factor intercept or total mean of endogenous factor is specified. Both cannot be simultaneously specified. |
| modelType | "CFA", "Sem", or "Path". This is specified to ensure that the analysis and data generation template created based on specified matrices in model correspond to what the user intends. |
| indLab | Character vector of indicator labels. If left blank, automatic labels will be generated as y1, y2, ... yy. |
| facLab | Character vector of factor labels. If left blank, automatic labels will be generated as f1, f2, ... ff |
| groupLab | Character of group-variable label (not the names of each group). If left blank, automatic labels will be generated as group |
| ngroups | Integer. Number of groups for data generation, defaults to 1. If larger than one, all specified matrices will be repeated for each additional group. If any matrix argument is a list, the length of this list will be the number of groups and ngroups is ignored. |
| smartStart | Defaults to FALSE. If TRUE, population parameter values that are real numbers will be used as starting values. |

Details

This function contains default settings:

For `modelType="CFA"`, LY is required. As the default, the on-diagonal elements of PS are fixed as 1 and the off-diagonal elements of PS are freely estimated. The off-diagonal elements of TE are freely estimated and the off-diagonal elements of TE are fixed to 0. The AL elements are fixed to 0. The TY elements are freely estimated.

For `modelType="Path"`, BE is required. As the default, the on-diagonal elements of PS are freely estimated, the off-diagonal elements between exogenous variables (covariance between exogenous variables) are freely estimated, and the other off-diagonal elements are fixed to 0. The AL elements are freely estimated.

For `modelType="SEM"`, LY and BE are required. As the default, the on-diagonal elements of PS are fixed to 1, the off-diagonal elements between exogenous factors (covariance between exogenous factors) are freely estimated, and the other off-diagonal elements are fixed to 0. The off-diagonal elements of TE are freely estimated and the off-diagonal elements of TE are fixed to 0. The AL elements are fixed to 0. The TY elements are freely estimated.

The `estmodel.cfa`, `estmodel.path`, and `estmodel.sem` are the shortcuts for the `estmodel` function when `modelType` are "CFA", "Path", and "SEM", respectively.

Value

SimSem object that contains the data generation template (@dgen) and analysis template (@pt).

Author(s)

Sunthud Pornprasertmanit (University of Kansas; <psunthud@ku.edu>)

See Also

- [model](#) To build data generation and data analysis template for simulation.
- [sim](#) For simulations using the [SimSem](#) template.
- [generate](#) To generate data using the [SimSem](#) template.
- [analyze](#) To analyze real or generated data using the [SimSem](#) template.
- [draw](#) To draw parameters using the [SimSem](#) template.

Examples

```
loading <- matrix(0, 6, 2)
loading[1:3, 1] <- NA
loading[4:6, 2] <- NA

CFA.Model <- estmodel(LY = loading, modelType = "CFA")
```

extractLavaanFit

Extract fit indices from the lavaan object

Description

Extract fit indices from the lavaan object

Usage

```
extractLavaanFit(Output)
```

Arguments

Output The lavaan object

Value

The renamed vector of fit measures

Author(s)

Sunthud Pornprasertmanit (University of Kansas; <psunthud@ku.edu>)

Examples

```
# No example
```

| | |
|------------|---|
| find2Dhist | <i>Fit the 2D Kernel Density Estimate</i> |
|------------|---|

Description

Fit the 2D Kernel Density Estimate to a pair of variables

Usage

```
find2Dhist(vec1, vec2)
```

Arguments

| | |
|------|------------|
| vec1 | Variable 1 |
| vec2 | Variable 2 |

Value

The 2D Kernel Density Estimate based on each pair of values in vec1 and vec2

Author(s)

Sunthud Pornprasertmanit (University of Kansas; <psunthud@ku.edu>)

Examples

```
# No example
```

| | |
|---------------------|--|
| findFactorIntercept | <i>Find factor intercept from regression coefficient matrix and factor total means</i> |
|---------------------|--|

Description

Find factor intercept from regression coefficient matrix and factor total means for latent variable models. In the path analysis model, this function will find indicator intercept from regression coefficient and indicator total means.

Usage

```
findFactorIntercept(beta, factorMean = NULL)
```

Arguments

| | |
|------------|---|
| beta | Regression coefficient matrix |
| factorMean | Total (model-implied) factor (indicator) means. The default is that all total factor means are 0. |

Value

A vector of factor (indicator) intercepts

Author(s)

Sunthud Pornprasertmanit (University of Kansas; <psunthud@ku.edu>)

See Also

- [findIndIntercept](#) to find indicator (measurement) intercepts
- [findIndMean](#) to find indicator (measurement) total means
- [findIndResidualVar](#) to find indicator (measurement) residual variances
- [findIndTotalVar](#) to find indicator (measurement) total variances
- [findFactorMean](#) to find factor means
- [findFactorResidualVar](#) to find factor residual variances
- [findFactorTotalVar](#) to find factor total variances
- [findFactorTotalCov](#) to find factor covariances

Examples

```
path <- matrix(0, 9, 9)
path[4, 1] <- path[7, 4] <- 0.6
path[5, 2] <- path[8, 5] <- 0.6
path[6, 3] <- path[9, 6] <- 0.6
path[5, 1] <- path[8, 4] <- 0.4
path[6, 2] <- path[9, 5] <- 0.4
factorMean <- c(5, 2, 3, 0, 0, 0, 0, 0, 0)
findFactorIntercept(path, factorMean)
```

| | |
|----------------|--|
| findFactorMean | <i>Find factor total means from regression coefficient matrix and factor intercept</i> |
|----------------|--|

Description

Find factor total means from regression coefficient matrix and factor intercepts for latent variable models. In the path analysis model, this function will find indicator total means from regression coefficient and indicator intercept.

Usage

```
findFactorMean(beta, alpha = NULL)
```

Arguments

| | |
|-------|--|
| beta | Regression coefficient matrix |
| alpha | Factor (indicator) intercept. The default is that all factor intercepts are 0. |

Value

A vector of factor (indicator) total means

Author(s)

Sunthud Pornprasertmanit (University of Kansas; <psunthud@ku.edu>)

See Also

- [findIndIntercept](#) to find indicator (measurement) intercepts
- [findIndMean](#) to find indicator (measurement) total means
- [findIndResidualVar](#) to find indicator (measurement) residual variances
- [findIndTotalVar](#) to find indicator (measurement) total variances
- [findFactorIntercept](#) to find factor intercepts
- [findFactorResidualVar](#) to find factor residual variances
- [findFactorTotalVar](#) to find factor total variances
- [findFactorTotalCov](#) to find factor covariances

Examples

```
path <- matrix(0, 9, 9)
path[4, 1] <- path[7, 4] <- 0.6
path[5, 2] <- path[8, 5] <- 0.6
path[6, 3] <- path[9, 6] <- 0.6
path[5, 1] <- path[8, 4] <- 0.4
path[6, 2] <- path[9, 5] <- 0.4
intcept <- c(5, 2, 3, 0, 0, 0, 0, 0, 0)
findFactorMean(path, intcept)
```

`findFactorResidualVar` *Find factor residual variances from regression coefficient matrix, factor (residual) correlations, and total factor variances*

Description

Find factor residual variances from regression coefficient matrix, factor (residual) correlation matrix, and total factor variances for latent variable models. In the path analysis model, this function will find indicator residual variances from regression coefficient, indicator (residual) correlation matrix, and total indicator variances.

Usage

```
findFactorResidualVar(beta, corPsi, totalVarPsi = NULL)
```

Arguments

| | |
|--------------------------|---|
| <code>beta</code> | Regression coefficient matrix |
| <code>corPsi</code> | Factor or indicator residual correlations. |
| <code>totalVarPsi</code> | Factor or indicator total variances. The default is that all factor or indicator total variances are 1. |

Value

A vector of factor (indicator) residual variances

Author(s)

Sunthud Pornprasertmanit (University of Kansas; <psunthud@ku.edu>)

See Also

- [findIndIntercept](#) to find indicator (measurement) intercepts
- [findIndMean](#) to find indicator (measurement) total means
- [findIndResidualVar](#) to find indicator (measurement) residual variances
- [findIndTotalVar](#) to find indicator (measurement) total variances
- [findFactorIntercept](#) to find factor intercepts
- [findFactorMean](#) to find factor means
- [findFactorTotalVar](#) to find factor total variances
- [findFactorTotalCov](#) to find factor covariances

Examples

```
path <- matrix(0, 9, 9)
path[4, 1] <- path[7, 4] <- 0.6
path[5, 2] <- path[8, 5] <- 0.6
path[6, 3] <- path[9, 6] <- 0.6
path[5, 1] <- path[8, 4] <- 0.4
path[6, 2] <- path[9, 5] <- 0.4
facCor <- diag(9)
facCor[1, 2] <- facCor[2, 1] <- 0.4
facCor[1, 3] <- facCor[3, 1] <- 0.4
facCor[2, 3] <- facCor[3, 2] <- 0.4
totalVar <- rep(1, 9)
findFactorResidualVar(path, facCor, totalVar)
```

| | |
|--------------------|--|
| findFactorTotalCov | <i>Find factor total covariance from regression coefficient matrix, factor residual covariance</i> |
|--------------------|--|

Description

Find factor total covariances from regression coefficient matrix, factor residual covariance matrix. The residual covariance matrix might be derived from factor residual correlation, total variance, and error variance. This function can be applied for path analysis model as well.

Usage

```
findFactorTotalCov(beta, psi=NULL, corPsi=NULL, totalVarPsi = NULL, errorVarPsi=NULL)
```

Arguments

| | |
|-------------|--|
| beta | Regression coefficient matrix |
| psi | Factor or indicator residual covariances. This argument can be skipped if factor residual correlation and either total variances or error variances are specified. |
| corPsi | Factor or indicator residual correlation. This argument must be specified with total variances or error variances. |
| totalVarPsi | Factor or indicator total variances. |
| errorVarPsi | Factor or indicator residual variances. |

Value

A matrix of factor (model-implied) total covariance

Author(s)

Sunthud Pornprasertmanit (University of Kansas; <psunthud@ku.edu>)

See Also

- [findIndIntercept](#) to find indicator (measurement) intercepts
- [findIndMean](#) to find indicator (measurement) total means
- [findIndResidualVar](#) to find indicator (measurement) residual variances
- [findIndTotalVar](#) to find indicator (measurement) total variances
- [findFactorIntercept](#) to find factor intercepts
- [findFactorMean](#) to find factor means
- [findFactorResidualVar](#) to find factor residual variances
- [findFactorTotalVar](#) to find factor total variances

Examples

```
path <- matrix(0, 9, 9)
path[4, 1] <- path[7, 4] <- 0.6
path[5, 2] <- path[8, 5] <- 0.6
path[6, 3] <- path[9, 6] <- 0.6
path[5, 1] <- path[8, 4] <- 0.4
path[6, 2] <- path[9, 5] <- 0.4
facCor <- diag(9)
facCor[1, 2] <- facCor[2, 1] <- 0.4
facCor[1, 3] <- facCor[3, 1] <- 0.4
facCor[2, 3] <- facCor[3, 2] <- 0.4
residualVar <- c(1, 1, 1, 0.64, 0.288, 0.288, 0.64, 0.29568, 0.21888)
findFactorTotalCov(path, corPsi=facCor, errorVarPsi=residualVar)
```

| | |
|--------------------|--|
| findFactorTotalVar | <i>Find factor total variances from regression coefficient matrix, factor (residual) correlations, and factor residual variances</i> |
|--------------------|--|

Description

Find factor total variances from regression coefficient matrix, factor (residual) correlation matrix, and factor residual variances for latent variable models. In the path analysis model, this function will find indicator total variances from regression coefficient, indicator (residual) correlation matrix, and indicator residual variances.

Usage

```
findFactorTotalVar(beta, corPsi, residualVarPsi)
```

Arguments

| | |
|----------------|--|
| beta | Regression coefficient matrix |
| corPsi | Factor or indicator residual correlations. |
| residualVarPsi | Factor or indicator residual variances. |

Value

A vector of factor (indicator) total variances

Author(s)

Sunthud Pornprasertmanit (University of Kansas; <psunthud@ku.edu>)

See Also

- [findIndIntercept](#) to find indicator (measurement) intercepts
- [findIndMean](#) to find indicator (measurement) total means
- [findIndResidualVar](#) to find indicator (measurement) residual variances
- [findIndTotalVar](#) to find indicator (measurement) total variances
- [findFactorIntercept](#) to find factor intercepts
- [findFactorMean](#) to find factor means
- [findFactorResidualVar](#) to find factor residual variances
- [findFactorTotalCov](#) to find factor covariances

Examples

```
path <- matrix(0, 9, 9)
path[4, 1] <- path[7, 4] <- 0.6
path[5, 2] <- path[8, 5] <- 0.6
path[6, 3] <- path[9, 6] <- 0.6
path[5, 1] <- path[8, 4] <- 0.4
path[6, 2] <- path[9, 5] <- 0.4
facCor <- diag(9)
facCor[1, 2] <- facCor[2, 1] <- 0.4
```

```

facCor[1, 3] <- facCor[3, 1] <- 0.4
facCor[2, 3] <- facCor[3, 2] <- 0.4
residualVar <- c(1, 1, 1, 0.64, 0.288, 0.288, 0.64, 0.29568, 0.21888)
findFactorTotalVar(path, facCor, residualVar)

```

| | |
|------------------|---|
| findIndIntercept | <i>Find indicator intercepts from factor loading matrix, total factor mean, and indicator mean.</i> |
|------------------|---|

Description

Find indicator (measurement) intercepts from a factor loading matrix, total factor mean, and indicator mean.

Usage

```
findIndIntercept(lambda, factorMean = NULL, indicatorMean = NULL)
```

Arguments

| | |
|---------------|--|
| lambda | Factor loading matrix |
| factorMean | Total (model-implied) mean of factors. As a default, all total factor means are 0. |
| indicatorMean | Total indicator means. As a default, all total indicator means are 0. |

Value

A vector of indicator (measurement) intercepts.

Author(s)

Sunthud Pornprasertmanit (University of Kansas; <psunthud@ku.edu>)

See Also

- [findIndMean](#) to find indicator (measurement) total means
- [findIndResidualVar](#) to find indicator (measurement) residual variances
- [findIndTotalVar](#) to find indicator (measurement) total variances
- [findFactorIntercept](#) to find factor intercepts
- [findFactorMean](#) to find factor means
- [findFactorResidualVar](#) to find factor residual variances
- [findFactorTotalVar](#) to find factor total variances
- [findFactorTotalCov](#) to find factor covariances

Examples

```

loading <- matrix(0, 6, 2)
loading[1:3, 1] <- c(0.6, 0.7, 0.8)
loading[4:6, 2] <- c(0.6, 0.7, 0.8)
facMean <- c(0.5, 0.2)
indMean <- rep(1, 6)
findIndIntercept(loading, facMean, indMean)

```

| | |
|-------------|---|
| findIndMean | <i>Find indicator total means from factor loading matrix, total factor mean, and indicator intercept.</i> |
|-------------|---|

Description

Find indicator total means from a factor loading matrix, total factor means, and indicator (measurement) intercepts.

Usage

```
findIndMean(lambda, factorMean = NULL, tau = NULL)
```

Arguments

| | |
|------------|--|
| lambda | Factor loading matrix |
| factorMean | Total (model-implied) mean of factors. As a default, all total factor means are 0. |
| tau | Indicator (measurement) intercepts. As a default, all intercepts are 0. |

Value

A vector of indicator total means.

Author(s)

Sunthud Pornprasertmanit (University of Kansas; <psunthud@ku.edu>)

See Also

- [findIndIntercept](#) to find indicator (measurement) intercepts
- [findIndResidualVar](#) to find indicator (measurement) residual variances
- [findIndTotalVar](#) to find indicator (measurement) total variances
- [findFactorIntercept](#) to find factor intercepts
- [findFactorMean](#) to find factor means
- [findFactorResidualVar](#) to find factor residual variances
- [findFactorTotalVar](#) to find factor total variances
- [findFactorTotalCov](#) to find factor covariances

Examples

```
loading <- matrix(0, 6, 2)
loading[1:3, 1] <- c(0.6, 0.7, 0.8)
loading[4:6, 2] <- c(0.6, 0.7, 0.8)
facMean <- c(0.5, 0.2)
intcept <- rep(0, 6)
findIndMean(loading, facMean, intcept)
```

| | |
|--------------------|--|
| findIndResidualVar | <i>Find indicator residual variances from factor loading matrix, total factor covariance, and total indicator variances.</i> |
|--------------------|--|

Description

Find indicator (measurement) residual variances from a factor loading matrix, total factor covariance matrix, and total indicator variances.

Usage

```
findIndResidualVar(lambda, totalFactorCov, totalVarTheta = NULL)
```

Arguments

| | |
|----------------|---|
| lambda | Factor loading matrix |
| totalFactorCov | Total (model-implied) covariance matrix among factors. |
| totalVarTheta | Indicator total variances. As a default, all total variances are 1. |

Value

A vector of indicator residual variances.

Author(s)

Sunthud Pornprasertmanit (University of Kansas; <psunthud@ku.edu>)

See Also

- [findIndIntercept](#) to find indicator (measurement) intercepts
- [findIndMean](#) to find indicator (measurement) total means
- [findIndTotalVar](#) to find indicator (measurement) total variances
- [findFactorIntercept](#) to find factor intercepts
- [findFactorMean](#) to find factor means
- [findFactorResidualVar](#) to find factor residual variances
- [findFactorTotalVar](#) to find factor total variances
- [findFactorTotalCov](#) to find factor covariances

Examples

```
loading <- matrix(0, 6, 2)
loading[1:3, 1] <- c(0.6, 0.7, 0.8)
loading[4:6, 2] <- c(0.6, 0.7, 0.8)
facCov <- matrix(c(1, 0.5, 0.5, 1), 2, 2)
totalVar <- rep(1, 6)
findIndResidualVar(loading, facCov, totalVar)
```

| | |
|-----------------|--|
| findIndTotalVar | <i>Find indicator total variances from factor loading matrix, total factor covariance, and indicator residual variances.</i> |
|-----------------|--|

Description

Find indicator total variances from a factor loading matrix, total factor covariance matrix, and indicator (measurement) residual variances.

Usage

```
findIndTotalVar(lambda, totalFactorCov, residualVarTheta)
```

Arguments

| | |
|------------------|--|
| lambda | Factor loading matrix |
| totalFactorCov | Total (model-implied) covariance matrix among factors. |
| residualVarTheta | Indicator (measurement) residual variances. |

Value

A vector of indicator total variances.

Author(s)

Sunthud Pornprasertmanit (University of Kansas; <psunthud@ku.edu>)

See Also

- [findIndIntercept](#) to find indicator (measurement) intercepts
- [findIndMean](#) to find indicator (measurement) total means
- [findIndResidualVar](#) to find indicator (measurement) residual variances
- [findFactorIntercept](#) to find factor intercepts
- [findFactorMean](#) to find factor means
- [findFactorResidualVar](#) to find factor residual variances
- [findFactorTotalVar](#) to find factor total variances
- [findFactorTotalCov](#) to find factor covariances

Examples

```
loading <- matrix(0, 6, 2)
loading[1:3, 1] <- c(0.6, 0.7, 0.8)
loading[4:6, 2] <- c(0.6, 0.7, 0.8)
facCov <- matrix(c(1, 0.5, 0.5, 1), 2, 2)
resVar <- c(0.64, 0.51, 0.36, 0.64, 0.51, 0.36)
findIndTotalVar(loading, facCov, resVar)
```

| | |
|-----------|--|
| findphist | <i>Find the density (likelihood) of a pair value in 2D Kernel Density Estimate</i> |
|-----------|--|

Description

Find the density (likelihood) of a pair value in 2D Kernel Density Estimate

Usage

```
findphist(value, hist)
```

Arguments

| | |
|-------|-------------------------------------|
| value | A target pair of values |
| hist | A 2D Binned Kernel Density Estimate |

Value

The probability (density) of the target pair of value

Author(s)

Sunthud Pornprasertmanit (University of Kansas; <psunthud@ku.edu>)

Examples

```
# No example
```

| | |
|-----------------------|---|
| findPossibleFactorCor | <i>Find the appropriate position for freely estimated correlation (or covariance) given a regression coefficient matrix</i> |
|-----------------------|---|

Description

Find the appropriate position for freely estimated correlation (or covariance) given a regression coefficient matrix. The appropriate position is the pair of variables that are not causally related.

Usage

```
findPossibleFactorCor(beta)
```

Arguments

| | |
|------|--|
| beta | The regression coefficient in path analysis. |
|------|--|

Value

The symmetric matrix containing the appropriate position for freely estimated correlation.

Author(s)

Sunthud Pornprasertmanit (University of Kansas; <psunthud@ku.edu>)

See Also

- [findRecursiveSet](#) to group variables regarding the position in mediation chain.

Examples

```
path <- matrix(0, 9, 9)
path[4, 1] <- path[7, 4] <- NA
path[5, 2] <- path[8, 5] <- NA
path[6, 3] <- path[9, 6] <- NA
path[5, 1] <- path[8, 4] <- NA
path[6, 2] <- path[9, 5] <- NA
findPossibleFactorCor(path)
```

findPower

Find a value of independent variables that provides a given value of power.

Description

Find a value of independent variable that provides a given value of power. If there are more than one varying parameters, this function will find the value of the target varying parameters given the values of the other varying parameters.

Usage

```
findPower(powerTable, iv, power)
```

Arguments

| | |
|------------|--|
| powerTable | A data.frame providing varying parameters and powers of each parameter. This table is obtained by getPower or continuousPower function. |
| iv | The target varying parameter that users would like to find the value providing a given power from. This argument can be specified as the index of the target column or the name of target column (i.e., "iv.N" or "N") |
| power | A desired power. |

Value

There are five possible types of values provided:

- *Value* The varying parameter value that provides the power just over the specified power value (the adjacent value of varying parameter provides lower power than the specified power value).
- *Minimum value* The minimum value has already provided enough power (way over the specified power value). The value of varying parameters that provides exact desired power may be lower than the minimum value. The example of varying parameter that can provides the minimum value is sample size.

- *Maximum value* The maximum value has already provided enough power (way over the specified power value). The value of varying parameters that provides exact desired power may be higher than the maximum value. The example of varying parameter that can provide the maximum value is percent missing.
- *NA* There is no value in the domain of varying parameters that provides the power greater than the desired power.
- *NaN* The power of all values in the varying parameters is 1 (specifically more than 0.9999) and any values of the varying parameters can be picked and still provide enough power.

Author(s)

Sunthud Pornprasertmanit (University of Kansas; <psunthud@ku.edu>)

See Also

- [getPower](#) to find the power of parameter estimates
- [continuousPower](#) to find the power of parameter estimates for the result object (`linkS4class{SimResult}`) with varying parameters.

Examples

```
## Not run:
# Specify Sample Size by n
loading <- matrix(0, 6, 1)
loading[1:6, 1] <- NA
LX <- bind(loading, 0.4)
RPH <- binds(diag(1))
RTD <- binds(diag(6))
CFA.Model <- model(LY = LX, RPS = RPH, RTE = RTD, modelType="CFA")

# Specify both sample size and percent missing completely at random
Output <- sim(NULL, model=CFA.Model, n=seq(100, 200, 20), pmMCAR=c(0, 0.1, 0.2))
pow <- getPower(Output)
findPower(pow, "N", 0.80)

## End(Not run)
```

findRecursiveSet

Group variables regarding the position in mediation chain

Description

In mediation analysis, variables affects other variables as a chain. This function will group variables regarding the chain of mediation analysis.

Usage

```
findRecursiveSet(beta)
```

Arguments

beta The regression coefficient in path analysis.

Value

The list of set of variables in the mediation chain. The variables in position 1 will be the independent variables. The variables in the last variables will be the end of the chain.

Author(s)

Sunthud Pornprasertmanit (University of Kansas; <psunthud@ku.edu>)

See Also

- [findPossibleFactorCor](#) to find the possible position for latent correlation given a regression coefficient matrix

Examples

```
path <- matrix(0, 9, 9)
path[4, 1] <- path[7, 4] <- NA
path[5, 2] <- path[8, 5] <- NA
path[6, 3] <- path[9, 6] <- NA
path[5, 1] <- path[8, 4] <- NA
path[6, 2] <- path[9, 5] <- NA
findRecursiveSet(path)
```

findRowZero

Find rows in a matrix that all elements are zero in non-fixed subset rows and columns.

Description

Find rows in a matrix that all elements are zero in non-fixed subset rows and columns. This function will be used in the [findRecursiveSet](#) function

Usage

```
findRowZero(square.matrix, is.row.fixed = FALSE)
```

Arguments

`square.matrix` Any square matrix
`is.row.fixed` A logical vector with the length equal to the dimension of the `square.matrix`. If TRUE, the function will skip examining this row.

Value

A vector of positions that contain rows of all zeros

Author(s)

Sunthud Pornprasertmanit (University of Kansas; <psunthud@ku.edu>)

Examples

```
# No example
```

| | |
|-----------------|---|
| findTargetPower | <i>Find a value of varying parameters that provides a given value of power.</i> |
|-----------------|---|

Description

Find a value of varying parameters that provides a given value of power. This function can deal with only one varying parameter only ([findPower](#) can deal with more than one varying parameter).

Usage

```
findTargetPower(iv, dv, power)
```

Arguments

| | |
|-------|---|
| iv | A vector of the target varying parameter |
| dv | A <code>data.frame</code> of the power table of target parameters |
| power | A desired power. |

Value

The value of the target varying parameter providing the desired power. If the value is NA, there is no value in the domain of varying parameters that provide the target power. If the value is the minimum value of the varying parameters, it means that the minimum value has already provided enough power. The value of varying parameters that provides exact desired power may be lower than the minimum value.

Author(s)

Sunthud Pornprasertmanit (University of Kansas; <psunthud@ku.edu>)

See Also

- [getPower](#) to find the power of parameter estimates
- [continuousPower](#) to find the power of parameter estimates for the result object (`linkS4class{SimResult}`) with varying parameters.
- [findPower](#) to find a value of varying parameters that provides a given value of power, which can deal with more than one varying parameter

Examples

```
# No example
```

| | |
|----------------|--|
| fitMeasuresChi | <i>Find fit indices from the discrepancy values of the target model and null models.</i> |
|----------------|--|

Description

Find fit indices from the discrepancy values of the target model and null models. This function is modified from the `fitMeasures` function in `lavaan` package

Usage

```
fitMeasuresChi(X2, df, p, X2.null, df.null, p.null, N, fit.measures="all")
```

Arguments

| | |
|--------------|---|
| X2 | The chi-square value of the target model |
| df | The degree of freedom of the target model |
| p | The p vlaue of the target model |
| X2.null | The chi-square value of the null model |
| df.null | The degree of freedom of the null model |
| p.null | The p value of the null model |
| N | Sample size |
| fit.measures | The list of selected fit measures |

Value

A vector of fit measures

Author(s)

Yves Rosseel in the `lavaan` package Modified by Sunthud Pornprasertmanit (University of Kansas; psunthud@ku.edu)

Examples

```
# No example
```

| | |
|----------|--|
| generate | <i>Generate data using SimSem template</i> |
|----------|--|

Description

This function can be used to generate random data based on the [SimSem](#) template. Some notable features include fine control of misspecification and misspecification optimization, as well as the ability to generate non-normal data. When using *simsem* for simulations, this function is used internally to generate data in the function `sim`, and can be helpful for debugging, or in creating data for use with other analysis programs.

Usage

```
generate(model, n, maxDraw=50, misfitBounds=NULL, misfitType="f0",
averageNumMisspec=FALSE, optMisfit=NULL, optDraws=50,
indDist=NULL, sequential=FALSE, facDist=NULL, errorDist=NULL,
indLab=NULL, modelBoot=FALSE, realData=NULL, params=FALSE)
```

Arguments

| | |
|--------------------------------|---|
| <code>model</code> | A SimSem object. |
| <code>n</code> | Integer of sample size. |
| <code>maxDraw</code> | Integer specifying the maximum number of attempts to draw a valid set of parameters (no negative error variance, standardized coefficients over 1). |
| <code>misfitBounds</code> | Vector that contains upper and lower bounds of the misfit measure. Sets of parameters drawn that are not within these bounds are rejected. |
| <code>misfitType</code> | Character vector indicating the fit measure used to assess the misfit of a set of parameters. Can be "f0", "rmsea", "srmr", or "all". |
| <code>averageNumMisspec</code> | TRUE or FALSE. ?? |
| <code>optMisfit</code> | Character vector of either "min" or "max" indicating either maximum or minimum optimized misfit. If not null, the set of parameters out of the number of draws in "optDraws" that has either the maximum or minimum misfit of the given misfit type will be returned. |
| <code>optDraws</code> | Number of parameter sets to draw if <code>optMisfit</code> is not null. The set of parameters with the maximum or minimum misfit will be returned. |
| <code>indDist</code> | A SimDataDist object or list of objects for a distribution of indicators. If one object is passed, each indicator will have the same distribution. Use when <code>sequential</code> is FALSE. |
| <code>sequential</code> | If TRUE, use a sequential method to create data such that the data from factor are generated first and apply to a set of equations to obtain the data of indicators. If FALSE, create data directly from model-implied mean and covariance of indicators. |
| <code>facDist</code> | A SimDataDist object or list of objects for the distribution of factors. If one object is passed, all factors will have the same distribution. Use when <code>sequential</code> is TRUE. |
| <code>errorDist</code> | An object or list of objects of type <code>SimDataDist</code> indicating the distribution of errors. If a single <code>SimDataDist</code> is specified, each error will be generated with that distribution. |
| <code>indLab</code> | A vector of indicator labels. When not specified, the variable names are <code>x1</code> , <code>x2</code> , ... <code>xN</code> . |
| <code>modelBoot</code> | When specified, a model-based bootstrap is used for data generation. See details for further information. This argument requires real data to be passed to <code>readData</code> . |
| <code>realData</code> | A <code>data.frame</code> containing real data. The data generated will follow the distribution of this data set. |
| <code>params</code> | If TRUE, return the parameters drawn along with the generated data set. Default is FALSE. |

Details

This function will use `mvnrm` function in MASS package to create data from model implied covariance matrix if the data distribution object (`SimDataDist`) is not specified. If the data distribution object is specified, the Gaussian copula model is used. See `SimDataDist` for further details. For the model-based bootstrap, the transformation proposed by Yung & Bentler (1996) is used. This procedure is the expansion from the Bollen and Stine (1992) bootstrap including a mean structure. The model-implied mean vector and covariance matrix with trivial misspecification will be used in the model-based bootstrap if `misspec` is specified. See page 133 of Bollen and Stine (1992) for a reference.

Internally, parameters are first drawn, and data is then created from these parameters. Both of these steps are available via the `draw` and `createData` functions respectively.

Value

A data.frame containing simulated data from the data generation template. A variable "group" is appended indicating group membership.

Author(s)

Sunthud Pornprasertmanit (University of Kansas; <psunthud@ku.edu>), Patrick Miller (University of Notre Dame; <pmille13@nd.edu>)

References

- Bollen, K. A., & Stine, R. A. (1992). Bootstrapping goodness-of-fit measures in structural equation models. *Sociological Methods and Research*, 21, 205-229.
- Yung, Y.-F., & Bentler, P. M. (1996). Bootstrapping techniques in analysis of mean and covariance structures. In G. A. Marcoulides & R. E. Schumacker (Eds.), *Advanced structural equation modeling: Issues and techniques* (pp. 195-226). Mahwah, NJ: Erlbaum.

See Also

- `draw` To draw parameters using the `SimSem` template.
- `createData` To generate random data using a set of parameters from `draw`

Examples

```
loading <- matrix(0, 6, 2)
loading[1:3, 1] <- NA
loading[4:6, 2] <- NA
LY <- bind(loading, 0.7)

latent.cor <- matrix(NA, 2, 2)
diag(latent.cor) <- 1
RPS <- binds(latent.cor, 0.5)

RTE <- binds(diag(6))

VY <- bind(rep(NA,6),2)

CFA.Model <- model(LY = LY, RPS = RPS, RTE = RTE, modelType = "CFA")

dat <- generate(CFA.Model,200)
```

| | |
|--------------|--|
| getCondQtile | <i>Get a quantile of a variable given values of predictors</i> |
|--------------|--|

Description

Find a quantile of a variable. If the predictors are specified, the result will provide the conditional quantile given specified value of predictors. The `quantreg` package is used to find conditional quantile.

Usage

```
getCondQtile(y, x=NULL, xval=NULL, df = 0, qtile = 0.5)
```

Arguments

| | |
|-------|--|
| y | The variable that users wish to find a quantile from |
| x | The predictors variables. If NULL, the unconditional quantile of the y is provided. |
| xval | The vector of predictors' values that users wish to find the conditional quantile from. If "all" is specified, the function will provide the conditional quantile of every value in x. |
| df | The degree of freedom used in spline method in predicting the fit indices by the predictors. If df is 0, the spline method will not be applied. |
| qtile | The quantile rank. |

Value

A (conditional) quantile value

Author(s)

Sunthud Pornprasertmanit (University of Kansas; <psunthud@ku.edu>)

See Also

- [getCutoff](#) for finding fit indices cutoffs using conditional quantiles

Examples

```
# No example
```

| | |
|-----------|---|
| getCutoff | <i>Find fit indices cutoff given a priori alpha level</i> |
|-----------|---|

Description

Extract fit indices information from the [SimResult](#) and getCutoff of fit indices given a priori alpha level

Usage

```
getCutoff(object, alpha, revDirec = FALSE, usedFit = NULL, ...)
```

Arguments

| | |
|----------|---|
| object | SimResult that saves the analysis results from multiple replications |
| alpha | A priori alpha level |
| revDirec | The default is to find critical point on the side that indicates worse fit (the right side of RMSEA or the left side of CFI). If specifying as TRUE, the directions are reversed. |
| usedFit | Vector of names of fit indices that researchers wish to getCutoffs from. The default is to getCutoffs of all fit indices. |
| ... | Additional arguments. |

Value

One-tailed cutoffs of several fit indices with a priori alpha level

Methods

signature(object="data.frame") This method will find the fit indices cutoff given a specified alpha level. The additional arguments are predictor, predictorVal, and df, which allows the fit indices predicted by any arbitrary independent variables (such as sample size or percent MCAR). The predictor is the data.frame of the predictor values. The number of rows of the predictor argument should be equal to the number of rows in the object. The predictorVal is the values of predictor that researchers would like to find the fit indices cutoffs from. The df is the degree of freedom used in spline method in predicting the fit indices by the predictors. If df is 0, the spline method will not be applied.

signature(object="matrix") The details are similar to the method for data.frame

signature(object="SimResult") This method will find the fit indices cutoff given a specified alpha level. The additional arguments are nVal, pmMCARval, pmMARval, and df, which are needed when using varying sample sizes or percent missing across replications in [SimResult](#). The nVal is the sample size value that researchers wish to find the fit indices cutoffs from. The pmMCARval is the percent missing completely at random value that researchers wish to find the fit indices cutoffs from. The pmMARval is the percent missing at random value that researchers wish to find the fit indices cutoffs from. The df is the degree of freedom used in spline method in predicting the fit indices by the predictors. If df is 0, the spline method will not be applied.

Author(s)

Sunthud Pornprasertmanit (University of Kansas; <psunthud@ku.edu>)

See Also

[SimResult](#) for a detail of simResult

Examples

```
## Not run:
loading <- matrix(0, 6, 2)
loading[1:3, 1] <- NA
loading[4:6, 2] <- NA
loadingValues <- matrix(0, 6, 2)
loadingValues[1:3, 1] <- 0.7
loadingValues[4:6, 2] <- 0.7
LX <- bind(loading, loadingValues)
latent.cor <- matrix(NA, 2, 2)
diag(latent.cor) <- 1
RPH <- binds(latent.cor, 0.5)
error.cor <- matrix(0, 6, 6)
diag(error.cor) <- 1
RTD <- binds(error.cor)
CFA.Model <- model(LY = LX, RPS = RPH, RTE = RTD, modelType="CFA")

# We make the examples running only 5 replications to save time.
# In reality, more replications are needed.
Output <- sim(5, n = 200, model=CFA.Model)
getCutoff(Output, 0.05)

# Finding the cutoff when the sample size is varied.
Output2 <- sim(NULL, model=CFA.Model, n=seq(50, 100, 10))
getCutoff(Output2, 0.05, nVal = 75)

## End(Not run)
```

| | |
|-----------------|---|
| getCutoffNested | <i>Find fit indices cutoff for nested model comparison given a priori alpha level</i> |
|-----------------|---|

Description

Extract fit indices information from the simulation of parent and nested models and getCutoff of fit indices given a priori alpha level

Usage

```
getCutoffNested(nested, parent, alpha = 0.05, usedFit = NULL, nVal = NULL, pmMCArval = NULL, pmM
```

Arguments

| | |
|--------|--|
| nested | SimResult that saves the analysis results of nested model from multiple replications |
| parent | SimResult that saves the analysis results of parent model from multiple replications |
| alpha | A priori alpha level |

| | |
|-----------|---|
| usedFit | Vector of names of fit indices that researchers wish to getCutoffs from. The default is to getCutoffs of all fit indices. |
| nVal | The sample size value that researchers wish to find the fit indices cutoffs from. |
| pmMCARval | The percent missing completely at random value that researchers wish to find the fit indices cutoffs from. |
| pmMARval | The percent missing at random value that researchers wish to find the fit indices cutoffs from. |
| df | The degree of freedom used in spline method in predicting the fit indices by the predictors. If df is 0, the spline method will not be applied. |

Value

One-tailed cutoffs of several fit indices with a priori alpha level

Author(s)

Sunthud Pornprasertmanit (University of Kansas; <psunthud@ku.edu>)

See Also

[SimResult](#) for a detail of simResult [getCutoff](#) for a detail of finding cutoffs for absolute fit

Examples

```
## Not run:
loading.null <- matrix(0, 6, 1)
loading.null[1:6, 1] <- NA
LX.NULL <- bind(loading.null, 0.7)
RPH.NULL <- binds(diag(1))

error.cor.mis <- matrix("rnorm(1, 0, 0.1)", 6, 6)
diag(error.cor.mis) <- 1
RTD <- binds(diag(6), misspec=error.cor.mis)
CFA.Model.NULL <- model(LY = LX.NULL, RPS = RPH.NULL, RTE = RTD, modelType="CFA")

loading.alt <- matrix(0, 6, 2)
loading.alt[1:3, 1] <- NA
loading.alt[4:6, 2] <- NA
LX.ALT <- bind(loading.alt, 0.7)
latent.cor.alt <- matrix(NA, 2, 2)
diag(latent.cor.alt) <- 1
RPH.ALT <- binds(latent.cor.alt, "runif(1, 0.7, 0.9)")
CFA.Model.ALT <- model(LY = LX.ALT, RPS = RPH.ALT, RTE = RTD, modelType="CFA")

# The actual number of replications should be greater than 10.
Output.NULL.NULL <- sim(10, n=500, model=CFA.Model.NULL, generate=CFA.Model.NULL)
Output.NULL.ALT <- sim(10, n=500, model=CFA.Model.ALT, generate=CFA.Model.NULL)

getCutoffNested(Output.NULL.NULL, Output.NULL.ALT)

## End(Not run)
```

| | |
|--------------------|---|
| getCutoffNonNested | <i>Find fit indices cutoff for non-nested model comparison given a priori alpha level</i> |
|--------------------|---|

Description

Extract fit indices information from the simulation of two models fitting on the datasets created from both models and getCutoff of fit indices given a priori alpha level

Usage

```
getCutoffNonNested(dat1Mod1, dat1Mod2, dat2Mod1=NULL, dat2Mod2=NULL,
alpha=.05, usedFit=NULL, onetailed=FALSE, nVal = NULL, pmMCARval = NULL,
pmMARval = NULL, df = 0)
```

Arguments

| | |
|-----------|---|
| dat1Mod1 | SimResult that saves the simulation of analyzing Model 1 by datasets created from Model 1 |
| dat1Mod2 | SimResult that saves the simulation of analyzing Model 2 by datasets created from Model 1 |
| dat2Mod1 | SimResult that saves the simulation of analyzing Model 1 by datasets created from Model 2 |
| dat2Mod2 | SimResult that saves the simulation of analyzing Model 2 by datasets created from Model 2 |
| alpha | A priori alpha level |
| usedFit | Vector of names of fit indices that researchers wish to get cutoffs from. The default is to get cutoffs of all fit indices. |
| onetailed | If TRUE, the function will find the cutoff from one-tail test. If FALSE, the function will find the cutoff from two-tailed test. |
| nVal | The sample size value that researchers wish to find the fit indices cutoffs from. |
| pmMCARval | The percent missing completely at random value that researchers wish to find the fit indices cutoffs from. |
| pmMARval | The percent missing at random value that researchers wish to find the fit indices cutoffs from. |
| df | The degree of freedom used in spline method in predicting the fit indices by the predictors. If df is 0, the spline method will not be applied. |

Value

One- or two-tailed cutoffs of several fit indices with a priori alpha level. The cutoff is based on the fit indices from Model 1 subtracted by the fit indices from Model 2.

Author(s)

Sunthud Pornprasertmanit (University of Kansas; <psunthud@ku.edu>)

See Also

[SimResult](#) for a detail of `simResult` [getCutoff](#) for a detail of finding cutoffs for absolute fit [getCutoffNested](#) for a detail of finding cutoffs for nested model comparison [plotCutoffNonNested](#) Plot cutoffs for non-nested model comparison

Examples

```
## Not run:
loading.A <- matrix(0, 8, 2)
loading.A[1:3, 1] <- NA
loading.A[4:8, 2] <- NA
LX.A <- bind(loading.A, 0.7)
latent.cor <- matrix(NA, 2, 2)
diag(latent.cor) <- 1
RPH <- binds(latent.cor, "runif(1, 0.7, 0.9)")
RTD <- binds(diag(8))
CFA.Model.A <- model(LY = LX.A, RPS = RPH, RTE = RTD, modelType="CFA")

loading.B <- matrix(0, 8, 2)
loading.B[1:4, 1] <- NA
loading.B[5:8, 2] <- NA
LX.B <- bind(loading.B, 0.7)
CFA.Model.B <- model(LY = LX.B, RPS = RPH, RTE = RTD, modelType="CFA")

# The actual number of replications should be greater than 10.
Output.A.A <- sim(10, n=500, model=CFA.Model.A, generate=CFA.Model.A)
Output.A.B <- sim(10, n=500, model=CFA.Model.B, generate=CFA.Model.A)
Output.B.A <- sim(10, n=500, model=CFA.Model.A, generate=CFA.Model.B)
Output.B.B <- sim(10, n=500, model=CFA.Model.B, generate=CFA.Model.B)

getCutoffNonNested(Output.A.A, Output.A.B, Output.B.A, Output.B.B)
getCutoffNonNested(Output.A.A, Output.A.B)
getCutoffNonNested(Output.B.B, Output.B.A)

## End(Not run)
```

getExtraOutput

*Get extra outputs from the result of simulation***Description**

Get extra outputs from a simulation result object ([SimResult](#)). Users can ask this package to extra output from the [lavaan](#) object in each iteration by setting the `outfun` argument (in the `sim` function). See the example below.

Usage

```
getExtraOutput(object)
```

Arguments

`object` [SimResult](#) that have the extra output extracted by the function defined in the `outfun` argument (in the `sim` function)

Value

A list of extra outputs

Author(s)

Sunthud Pornprasertmanit (University of Kansas; <psunthud@ku.edu>)

See Also

- [sim](#) A function to run a Monte Carlo simulation

Examples

```
## Not run:
# Specify Sample Size by n
loading <- matrix(0, 6, 1)
loading[1:6, 1] <- NA
LX <- bind(loading, 0.7)
RPH <- binds(diag(1))
RTD <- binds(diag(6))
CFA.Model <- model(LY = LX, RPS = RPH, RTE = RTD, modelType="CFA")

# We will use only 5 replications to save time.
# In reality, more replications are needed.

outfun <- function(out) {
  result <- inspect(out, "mi")
}

# Specify both sample size and percent missing completely at random
Output <- sim(5, n=200, model=CFA.Model, outfun=outfun)
getExtraOutput(Output)

## End(Not run)
```

getKeywords

List of all keywords used in the simsem package

Description

List of all keywords used in the simsem package

Usage

```
getKeywords()
```

Value

A list of all keywords used in this package

- `usedFit` Fit indices used as the default for providing output
- `usedFitPop` Population fit indices used as the default for providing input
- `optMin` The method picking the minimum value of misfit across misspecification sets

- optMax The method picking the maximum value of misfit across misspecification sets
- optNone Not using the optimization method

Author(s)

Sunthud Pornprasertmanit (University of Kansas; psunthud@ku.edu)

Examples

```
# This function is not a public function.

# getKeywords()
```

| | |
|---------------|--|
| getPopulation | <i>Extract the data generation population model underlying a result object</i> |
|---------------|--|

Description

This function will extract the data generation population model underlying a result object (`linkS4class{SimResult}`).

Usage

```
getPopulation(object)
```

Arguments

| | |
|--------|---|
| object | The result object that you wish to extract the data generation population model from (<code>linkS4class{SimResult}</code>). |
|--------|---|

Value

A data frame contained the population of each replication

Author(s)

Sunthud Pornprasertmanit (University of Kansas; <psunthud@ku.edu>)

See Also

- [SimResult](#) for result object

Examples

```
## Not run:
loading <- matrix(0, 6, 1)
loading[1:6, 1] <- NA
LX <- bind(loading, "runif(1, 0.4, 0.9)")
RPH <- binds(diag(1))
RTD <- binds(diag(6))
CFA.Model <- model(LY = LX, RPS = RPH, RTE = RTD, modelType="CFA")

# We will use only 10 replications to save time.
```

```
# In reality, more replications are needed.
Output <- sim(10, n=200, model=CFA.Model)
getPopulation(Output)

## End(Not run)
```

getPower

Find power of model parameters

Description

A function to find the power of parameters in a model when none, one, or more of the simulations parameters vary randomly across replications.

Usage

```
getPower(simResult, alpha = 0.05, contParam = NULL, powerParam = NULL,
nVal = NULL, pmMCARval = NULL, pmMARval = NULL, paramVal = NULL)
```

Arguments

| | |
|------------|--|
| simResult | SimResult that may include at least one randomly varying parameter (e.g. sample size, percent missing, model parameters) |
| alpha | Alpha level to use for power analysis. |
| contParam | Vector of parameters names that vary over replications. |
| powerParam | Vector of parameters names that the user wishes to find power for. This can be a vector of names (e.g., "LY1_1", "LY2_2"), or the name of a matrix (e.g. "PS"), if the name of a matrix is used power for all parameters in that matrix will be returned. If parameters are not specified, power for all parameters in the model will be returned. |
| nVal | The sample size values that users wish to find power from. |
| pmMCARval | The percent completely missing at random values that users wish to find power from. |
| pmMARval | The percent missing at random values that users wish to find power from. |
| paramVal | A list of varying parameter values that users wish to find power from. |

Details

A common use of simulations is to conduct power analyses, especially when using SEM (Muthen & Muthen, 2002). Here, researchers could select values for each parameter and a sample size and run a simulation to determine power in those conditions (the proportion of generated datasets in which a particular parameter of interest is significantly different from zero). To evaluate power at multiple sample sizes, one simulation for each sample size must be run. This function not only calculate power for each sample size but also calculate power for multiple sample sizes varying continuously. By continuously varying sample size across replications, only a single simulation is needed. In this simulation, the sample size for each replication varies randomly across plausible sample sizes (e.g., sample sizes between 200 and 500). For each replication, the sample size and significance of each parameter (0 = not significant, 1 = significant) are recorded. When the simulation is complete, parameter significance is regressed on sample size using logistic regression. For a given sample size, the predicted probability from the logistic regression equation is the power to detect an effect at that sample size. This approach can be extended to other randomly varying simulation parameters such as the percentage of missing data, and model parameters.

Value

Data frame containing columns representing values of the randomly varying simulation parameters, and power for model parameters of interest.

Author(s)

Alexander M. Schoemann (University of Kansas; <schoemann@ku.edu>), Sunthud Pornprasertmanit (University of Kansas; <psunthud@ku.edu>)

References

Muthen, L. K., & Muthen, B. O. (2002). How to use a Monte Carlo study to decide on sample size and determine power. *Structural Equation Modeling*, 4, 599-620.

See Also

- [SimResult](#) to see how to create a simResult object with randomly varying parameters.

Examples

```
## Not run:
# Specify Sample Size by n
loading <- matrix(0, 6, 1)
loading[1:6, 1] <- NA
LX <- bind(loading, 0.7)
RPH <- binds(diag(1))
RTD <- binds(diag(6))
CFA.Model <- model(LY = LX, RPS = RPH, RTE = RTD, modelType="CFA")

# We will use only 5 replications to save time.
# In reality, more replications are needed.

# Specify both sample size and percent missing completely at random
Output <- sim(NULL, model=CFA.Model, n=seq(100, 200, 20), pmMCAR=c(0, 0.1, 0.2))
summary(Output)

getPower(Output)

getPower(Output, nVal=c(100, 200), pmMCARval=c(0, 0.1, 0.2))

## End(Not run)
```

getPowerFit

Find power in rejecting alternative models based on fit indices criteria

Description

Find the proportion of fit indices that indicate worse fit than a specified cutoffs. The cutoffs may be calculated from [getCutoff](#) of the null model.

Usage

```
getPowerFit(altObject, cutoff, revDirec = FALSE, usedFit=NULL, ...)
```

Arguments

| | |
|-----------|--|
| altObject | SimResult that indicates alternative model that users wish to reject |
| cutoff | Fit indices cutoffs from null model or users. This should be a vector with a specified fit indices names as the name of vector elements. This argument can be missing if the SimResult is specified in the altObject and the SimResult of the null model is specified. |
| revDirec | The default is to count the proportion of fit indices that indicates lower fit to the model, such as how many RMSEA in the alternative model that is worse than cutoffs. The direction can be reversed by setting as TRUE. |
| usedFit | Vector of names of fit indices that researchers wish to getCutoffs from. The default is to getCutoffs of all fit indices. |
| ... | Additional arguments |

Value

List of power given different fit indices.

Methods

signature(altObject="data.frame", cutoff="vector") This method will find the fit indices indicated in the altObject that provides worse fit than the cutoff. The additional arguments are predictor, predictorVal, condCutoff, and df, which allows the fit indices predicted by any arbitrary independent variables (such as sample size or percent MCAR). The predictor is the data.frame of the predictor values. The number of rows of the predictor argument should be equal to the number of rows in the object. The predictorVal is the values of predictor that researchers would like to find the power from. The condCutoff is a logical. If TRUE, the cutoff is applicable only a given value of predictorVal. If FALSE, the cutoff is applicable in any values of predictor. The df is the degree of freedom used in spline method in predicting the fit indices by the predictors. If df is 0, the spline method will not be applied.

signature(altObject="matrix", cutoff="vector") The details are similar to the method for altObject="data.frame" and cutoff="vector".

signature(altObject="SimResult", cutoff="vector") This method will find the fit indices indicated in the altObject that provides worse fit than the cutoff. The additional arguments are nVal, pmMCARval, pmMARval, condCutoff, and df, which are needed when using varying sample sizes or percent missing across replications in [SimResult](#). The nVal is the sample size value that researchers wish to find the fit indices cutoffs from. The pmMCARval is the percent missing completely at random value that researchers wish to find the fit indices cutoffs from. The pmMARval is the percent missing at random value that researchers wish to find the fit indices cutoffs from. The condCutoff is a logical. If TRUE, the cutoff is applicable only a given set of nVal, pmMCARval, and pmMARval. If FALSE, the cutoff is applicable in any values of sample size and percent missing. The df is the degree of freedom used in spline method in predicting the fit indices by the predictors. If df is 0, the spline method will not be applied.

signature(altObject="SimResult", cutoff="missing") The details are similar to the method for altObject="SimResult" and cutoff="vector". The cutoff argument must not be specified. Rather, the nullObject, which is an additional argument of this method, is required. The nullObject is the [SimResult](#) that contains the simulation result from fitting the null model by the data from the null model.

Author(s)

Sunthud Pornprasertmanit (University of Kansas; <psunthud@ku.edu>)

See Also

- [getCutoff](#) to find the cutoffs from null model.
- [SimResult](#) to see how to create simResult

Examples

```
## Not run:
loading.null <- matrix(0, 6, 1)
loading.null[1:6, 1] <- NA
LX.NULL <- bind(loading.null, 0.7)
RPH.NULL <- binds(diag(1))
RTD <- binds(diag(6))
CFA.Model.NULL <- model(LY = LX.NULL, RPS = RPH.NULL, RTE = RTD, modelType="CFA")

# We make the examples running only 5 replications to save time.
# In reality, more replications are needed.
Output.NULL <- sim(5, n=500, model=CFA.Model.NULL)
Cut.NULL <- getCutoff(Output.NULL, 0.95)

loading.alt <- matrix(0, 6, 2)
loading.alt[1:3, 1] <- NA
loading.alt[4:6, 2] <- NA
LX.ALT <- bind(loading.alt, 0.7)
latent.cor.alt <- matrix(NA, 2, 2)
diag(latent.cor.alt) <- 1
RPH.ALT <- binds(latent.cor.alt, "runif(1, 0.7, 0.9)")
CFA.Model.ALT <- model(LY = LX.ALT, RPS = RPH.ALT, RTE = RTD, modelType="CFA")

Output.ALT <- sim(5, n=500, model=CFA.Model.NULL, generate=CFA.Model.ALT)
getPowerFit(Output.ALT, cutoff=Cut.NULL)
Rule.of.thumb <- c(RMSEA=0.05, CFI=0.95, TLI=0.95, SRMR=0.06)
getPowerFit(Output.ALT, cutoff=Rule.of.thumb, usedFit=c("RMSEA", "CFI", "TLI", "SRMR"))

Output.NULL2 <- sim(NULL, n=seq(50, 500, 50), model=CFA.Model.NULL, generate=CFA.Model.NULL)
Output.ALT2 <- sim(NULL, n=seq(50, 500, 50), model=CFA.Model.NULL, generate=CFA.Model.ALT)
getPowerFit(Output.ALT2, nullObject=Output.NULL2, nVal=250)

## End(Not run)
```

| | |
|-------------------|--|
| getPowerFitNested | <i>Find power in rejecting nested models based on the differences in fit indices</i> |
|-------------------|--|

Description

Find the proportion of the difference in fit indices that indicate worse fit than a specified (or internally derived) cutoffs.

Usage

```
getPowerFitNested(altNested, altParent, cutoff, ...)
```

Arguments

| | |
|-----------|--|
| altNested | SimResult that saves the simulation result of the nested model when the nested model is FALSE. |
| altParent | SimResult that saves the simulation result of the parent model when the nested model is FALSE. |
| cutoff | A vector of priori cutoffs for fit indices. |
| ... | Additional arguments |

Value

List of power given different fit indices.

Methods

signature(altNested="SimResult", altParent="SimResult", cutoff="vector") This method will find the the differences in fit indices from altNested and altParent that provides worse fit than the cutoff. The additional arguments are revDirec, usedFit, nVal, pmMCARval, pmMARval, condCutoff, and df, which are needed when using varying sample sizes or percent missing across replications in [SimResult](#). The revDirec is whether to reverse a direction. The default is to count the proportion of fit indices that indicates lower fit to the model, such as how many RMSEA in the alternative model that is worse than cutoffs. The direction can be reversed by setting as TRUE. The usedFit is the vector of names of fit indices that researchers wish to get power from. The default is to get the powers of all fit indices. The nVal is the sample size value that researchers wish to find the fit indices cutoffs from. The pmMCARval is the percent missing completely at random value that researchers wish to find the fit indices cutoffs from. The pmMARval is the percent missing at random value that researchers wish to find the fit indices cutoffs from. The condCutoff is a logical. If TRUE, the cutoff is applicable only a given set of nVal, pmMCARval, and pmMARval. If FALSE, the cutoff is applicable in any values of sample size and percent missing. The df is the degree of freedom used in spline method in predicting the fit indices by the predictors. If df is 0, the spline method will not be applied.

signature(altNested="SimResult", altParent="SimResult", cutoff="missing") The details are similar to the method for altNested="SimResult", altParent="SimResult", and cutoff="vector". The cutoff argument must not be specified. Rather, the nullNested and nullParent, which are additional arguments of this method, are required. The nullNested is the [SimResult](#) that saves the simulation result of the nested model when the nested model is TRUE. The nullParent is the [SimResult](#) that saves the simulation result of the parent model when the nested model is TRUE.

Author(s)

Sunthud Pornprasertmanit (University of Kansas; <psunthud@ku.edu>)

See Also

- [getCutoff](#) to find the cutoffs from null model.
- [SimResult](#) to see how to create simResult

Examples

```
## Not run:
loading.null <- matrix(0, 6, 1)
loading.null[1:6, 1] <- NA
LX.NULL <- bind(loading.null, 0.7)
RPH.NULL <- binds(diag(1))
RTD <- binds(diag(6))
CFA.Model.NULL <- model(LY = LX.NULL, RPS = RPH.NULL, RTE = RTD, modelType="CFA")

loading.alt <- matrix(0, 6, 2)
loading.alt[1:3, 1] <- NA
loading.alt[4:6, 2] <- NA
LX.ALT <- bind(loading.alt, 0.7)
latent.cor.alt <- matrix(NA, 2, 2)
diag(latent.cor.alt) <- 1
RPH.ALT <- binds(latent.cor.alt, 0.7)
CFA.Model.ALT <- model(LY = LX.ALT, RPS = RPH.ALT, RTE = RTD, modelType="CFA")

Output.NULL.NULL <- sim(10, n=500, model=CFA.Model.NULL, generate=CFA.Model.NULL)
Output.ALT.NULL <- sim(10, n=500, model=CFA.Model.NULL, generate=CFA.Model.ALT)
Output.NULL.ALT <- sim(10, n=500, model=CFA.Model.ALT, generate=CFA.Model.NULL)
Output.ALT.ALT <- sim(10, n=500, model=CFA.Model.ALT, generate=CFA.Model.ALT)

getPowerFitNested(Output.ALT.NULL, Output.ALT.ALT, nullNested=Output.NULL.NULL, nullParent=Output.NULL.ALT)
getPowerFitNested(Output.ALT.NULL, Output.ALT.ALT, cutoff=c(Chi=3.84, CFI=-0.10))

Output.NULL.NULL2 <- sim(NULL, n=seq(50, 500, 50), model=CFA.Model.NULL, generate=CFA.Model.NULL)
Output.ALT.NULL2 <- sim(NULL, n=seq(50, 500, 50), model=CFA.Model.NULL, generate=CFA.Model.ALT)
Output.NULL.ALT2 <- sim(NULL, n=seq(50, 500, 50), model=CFA.Model.ALT, generate=CFA.Model.NULL)
Output.ALT.ALT2 <- sim(NULL, n=seq(50, 500, 50), model=CFA.Model.ALT, generate=CFA.Model.ALT)

getPowerFitNested(Output.ALT.NULL2, Output.ALT.ALT2, nullNested=Output.NULL.NULL2, nullParent=Output.NULL.ALT2)
getPowerFitNested(Output.ALT.NULL2, Output.ALT.ALT2, cutoff=c(Chi=3.84, CFI=-0.10), nVal = 250)

## End(Not run)
```

| | |
|----------------------|--|
| getPowerFitNonNested | <i>Find power in rejecting non-nested models based on the differences in fit indices</i> |
|----------------------|--|

Description

Find the proportion of the difference in fit indices from one model that does not in the range of sampling distribution from another model (reject that the dataset comes from the second model) or indicates worse fit than a specified cutoff.

Usage

```
getPowerFitNonNested(dat2Mod1, dat2Mod2, cutoff, ...)
```

Arguments

| | |
|----------|---|
| dat2Mod1 | SimResult that saves the simulation of analyzing Model 1 by datasets created from Model 2 |
|----------|---|

| | |
|----------|---|
| dat2Mod2 | SimResult that saves the simulation of analyzing Model 2 by datasets created from Model 2 |
| cutoff | A vector of priori cutoffs for fit indices. |
| ... | Additional arguments |

Value

List of power given different fit indices.

Methods

signature(dat2Mod1="SimResult", dat2Mod2="SimResult", cutoff="vector") This method will find the the differences in fit indices from dat2Mod1 and dat2Mod2 that provides worse fit than the cutoff. The additional arguments are revDirec, usedFit, nVal, pmMCARval, pmMARval, condCutoff, and df, which are needed when using varying sample sizes or percent missing across replications in [SimResult](#). The revDirec is whether to reverse a direction. The default is to count the proportion of the difference of fit indices that lower than the specified cutoffs, such as how many the difference in RMSEA in the alternative model that is lower than cutoffs. The direction can be reversed by setting as TRUE. The usedFit is the vector of names of fit indices that researchers wish to get power from. The default is to get the powers of all fit indices. The nVal is the sample size value that researchers wish to find the fit indices cutoffs from. The pmMCARval is the percent missing completely at random value that researchers wish to find the fit indices cutoffs from. The pmMARval is the percent missing at random value that researchers wish to find the fit indices cutoffs from. The condCutoff is a logical. If TRUE, the cutoff is applicable only a given set of nVal, pmMCARval, and pmMARval. If FALSE, the cutoff is applicable in any values of sample size and percent missing. The df is the degree of freedom used in spline method in predicting the fit indices by the predictors. If df is 0, the spline method will not be applied.

signature(dat2Mod1="SimResult", dat2Mod2="SimResult", cutoff="missing") The details are similar to the method for dat2Mod1="SimResult", dat2Mod2="SimResult", and cutoff="vector". The cutoff argument must not be specified. Rather, the dat1Mod1 and dat1Mod2, which are additional arguments of this method, are required. The dat1Mod1 is the [SimResult](#) that saves the simulation of analyzing Model 1 by datasets created from Model 1. The dat1Mod2 is the [SimResult](#) that saves the simulation of analyzing Model 2 by datasets created from Model 1. The another additional argument is onetailed that is to derive the cutoff by using one-tailed test if specified as TRUE.

Author(s)

Sunthud Pornprasertmanit (University of Kansas; <psunthud@ku.edu>)

See Also

- [getCutoffNonNested](#) to find the cutoffs for non-nested model comparison
- [SimResult](#) to see how to create simResult

Examples

```
## Not run:
loading.A <- matrix(0, 8, 2)
loading.A[1:3, 1] <- NA
loading.A[4:8, 2] <- NA
LX.A <- bind(loading.A, 0.7)
```

```

latent.cor <- matrix(NA, 2, 2)
diag(latent.cor) <- 1
RPH <- binds(latent.cor, "runif(1, 0.7, 0.9)")
RTD <- binds(diag(8))
CFA.Model.A <- model(LY = LX.A, RPS = RPH, RTE = RTD, modelType="CFA")

loading.B <- matrix(0, 8, 2)
loading.B[1:4, 1] <- NA
loading.B[5:8, 2] <- NA
LX.B <- bind(loading.B, 0.7)
CFA.Model.B <- model(LY = LX.B, RPS = RPH, RTE = RTD, modelType="CFA")

# The actual number of replications should be greater than 10.
Output.A.A <- sim(10, n=500, model=CFA.Model.A, generate=CFA.Model.A)
Output.A.B <- sim(10, n=500, model=CFA.Model.B, generate=CFA.Model.A)
Output.B.A <- sim(10, n=500, model=CFA.Model.A, generate=CFA.Model.B)
Output.B.B <- sim(10, n=500, model=CFA.Model.B, generate=CFA.Model.B)

getPowerFitNonNested(Output.B.A, Output.B.B, dat1Mod1=Output.A.A, dat1Mod2=Output.A.B)
getPowerFitNonNested(Output.B.A, Output.B.B, cutoff=c(AIC=0, BIC=0))

## End(Not run)

```

imposeMissing

Impose MAR, MCAR, planned missingness, or attrition on a data set

Description

Function imposes missing values on a data based on the known missing data types, including MCAR, MAR, planned, and attrition.

Usage

```

impose(miss, data.mat, pmMCAR = NULL, pmMAR = NULL)
imposeMissing(data.mat, cov = 0, pmMCAR = 0, pmMAR = 0, nforms = 0,
itemGroups = list(), twoMethod = 0, prAttr = 0, timePoints = 1,
ignoreCols = 0, threshold = 0, logical = NULL)

```

Arguments

| | |
|----------|---|
| miss | Missing data object (SimMissing) used as the template for impose missing values |
| data.mat | Data to impose missing upon. Can be either a matrix or a data frame. |
| cov | Column indices of a covariate to be used to impose MAR missing, or MAR attrition. Will not be included in any removal procedure. See details. |
| pmMCAR | Decimal percent of missingness to introduce completely at random on all variables. |
| pmMAR | Decimal percent of missingness to introduce using the listed covariate as predictor. See details. |
| nforms | The number of forms for planned missing data designs, not including the shared form. |

| | |
|-------------------------|---|
| <code>itemGroups</code> | List of lists of item groupings for planned missing data forms. Unless specified, items will be divided into groups sequentially (e.g. 1-3,4-6,7-9,10-12) |
| <code>twoMethod</code> | With missing on one variable: vector of (column index, percent missing). Will put a given percent missing on that column in the matrix to simulate a two method planned missing data research design. With missing on two or more variables: list of (column indices, percent missing). |
| <code>prAttr</code> | Probability (or vector of probabilities) of an entire case being removed due to attrition at a given time point. When a covariate is specified along with this argument, attrition will be predicted by the covariate (MAR attrition). See details. |
| <code>timePoints</code> | Number of timepoints items were measured over. For longitudinal data, planned missing designs will be implemented within each timepoint. All methods to impose missing values over time assume an equal number of variables at each time point. |
| <code>ignoreCols</code> | The columns not imposed any missing values for any missing data patterns. |
| <code>threshold</code> | The threshold of the covariate used to impose missing values. Values on the covariate above this threshold are eligible to be deleted. The default threshold is the mean of the variable. |
| <code>logical</code> | A matrix of logical values (TRUE/FALSE). If a value in the dataset is corresponding to the TRUE in the logical matrix, the value will be missing. |

Details

Without specifying any arguments, no missing values will be introduced.

A single covariate is required to specify MAR missing - this covariate can be distributed in any way. This covariate can be either continuous or categorical, as long as it is numerical. If the covariate is categorical, the threshold should be specified to one of the levels.

MAR missingness is specified using the threshold method - any value on the covariate that is above the specified threshold indicates a row eligible for deletion. If the specified total amount of MAR missingness is not possible given the total rows eligible based on the threshold, the function iteratively lowers the threshold until the total percent missing is possible.

Planned missingness is parameterized by the number of forms (n). This is used to divide the cases into n groups. If the column groupings are not specified, a naive method will be used that divides the columns into $n+1$ equal forms sequentially (1-4,5-9,10-13..), where the first group is the shared form. The first list of column indices given will be used as the shared group. If this is not desired, this list can be left empty.

For attrition, the probability can be specified as a single value or as a vector. For a single value, the probability of attrition will be the same across time points, and affects only cases not previously lost due to attrition. If this argument is a vector, this specifies different probabilities of attrition for each time point. Values will be recycled if this vector is smaller than the specified number of time points.

An MNAR processes can be generated by specifying MAR missingness and then dropping the covariate from the subsequent analysis.

Currently, if MAR missing is imposed along with attrition, both processes will use the same covariate and threshold.

Currently, all types of missingness (MCAR, MAR, planned, and attrition) are imposed independently. This means that specified global values of percent missing will not be additive (10 percent MCAR + 10 percent MAR does not equal 20 percent total missing).

Value

A data matrix with NAs introduced in the way specified by the arguments.

Author(s)

Patrick Miller(University of Kansas; <patr1ckm@ku.edu>) Alexander M. Schoemann (University of Kansas; <schoemann@ku.edu>)

See Also

- [SimMissing](#) for the alternative way to save missing data feature for using in the [sim](#) function.

Examples

```
data <- matrix(rep(rnorm(10,1,1),19),ncol=19)
datac <- cbind(data,rnorm(10,0,1),rnorm(10,5,5))

# Imposing Missing with the following arguments produces no missing values
imposeMissing(data)
imposeMissing(data,cov=c(1,2))
imposeMissing(data,pmMCAR=0)
imposeMissing(data,pmMAR=0)
imposeMissing(data,nforms=0)

#Some more usage examples
imposeMissing(data,cov=c(1,2),pmMCAR=.1)

imposeMissing(data,nforms=3)
imposeMissing(data,nforms=3,itemGroups=list(c(1,2,3,4,5),c(6,7,8,9,10),c(11,12,13,14,15),c(16,17,18,19)))
imposeMissing(datac,cov=c(20,21),nforms=3)
imposeMissing(data,twoMethod=c(19,.8))
imposeMissing(datac,cov=21,prAttr=.1,timePoints=5)
```

interpolate

Find the value of one vector relative to a value of another vector by interpolation

Description

Find the value of the resulting vector that have the position similar to the value of the baseline vector. If the starting value in the baseline vector is in between two elements, the resulting value will be predicted by linear interpolation.

Usage

```
interpolate(baselineVec, val, resultVec=NULL)
```

Arguments

| | |
|-------------|---|
| baselineVec | The target vector to be used as a baseline. The resulting vector can be attached as the element names of this vector. |
| val | The value relative to the baseline vector to be used for projecting the resulting value |
| resultVec | The vector that the resulting value will be used to base their result form |

Value

The interpolated value from the resulting vector relative to the value in the baseline vector

Author(s)

Sunthud Pornprasertmanit (University of Kansas; <psunthud@ku.edu>)

Examples

```
# No Example
```

| | |
|-------------|---|
| likRatioFit | <i>Find the likelihood ratio (or Bayes factor) based on the bivariate distribution of fit indices</i> |
|-------------|---|

Description

Find the log-likelihood of the observed fit indices on Model 1 and 2 from the real data on the bivariate sampling distribution of fit indices fitting Model 1 and Model 2 by the datasets from the Model 1 and Model 2. Then, the likelihood ratio is computed (which may be interpreted as posterior odd). If the prior odd is 1 (by default), the likelihood ratio is equivalent to Bayes Factor.

Usage

```
likRatioFit(outMod1, outMod2, dat1Mod1, dat1Mod2, dat2Mod1, dat2Mod2,
  usedFit=NULL, prior=1)
```

Arguments

| | |
|----------|---|
| outMod1 | lavaan that saves the analysis result of the first model from the target dataset |
| outMod2 | lavaan that saves the analysis result of the second model from the target dataset |
| dat1Mod1 | SimResult that saves the simulation of analyzing Model 1 by datasets created from Model 1 |
| dat1Mod2 | SimResult that saves the simulation of analyzing Model 2 by datasets created from Model 1 |
| dat2Mod1 | SimResult that saves the simulation of analyzing Model 1 by datasets created from Model 2 |
| dat2Mod2 | SimResult that saves the simulation of analyzing Model 2 by datasets created from Model 2 |
| usedFit | Vector of names of fit indices that researchers wish to getCutoffs from. The default is to getCutoffs of all fit indices. |
| prior | The prior odds. The prior probability that Model 1 is correct over the prior probability that Model 2 is correct. |

Value

The likelihood ratio (Bayes Factor) in preference of Model 1 to Model 2. If the value is greater than 1, Model 1 is preferred. If the value is less than 1, Model 2 is preferred.

Author(s)

Sunthud Pornprasertmanit (University of Kansas; <psunthud@ku.edu>)

See Also

[SimResult](#) for a detail of `simResult` [pValueNested](#) for a nested model comparison by the difference in fit indices [pValueNonNested](#) for a nonnested model comparison by the difference in fit indices

Examples

```
## Not run:
library(lavaan)
loading <- matrix(0, 11, 3)
loading[1:3, 1] <- NA
loading[4:7, 2] <- NA
loading[8:11, 3] <- NA
path.A <- matrix(0, 3, 3)
path.A[2:3, 1] <- NA
path.A[3, 2] <- NA
model.A <- estmodel(LY=loading, BE=path.A, modelType="SEM", indLab=c(paste("x", 1:3, sep=""), paste("y", 1:3, sep="")))

out.A <- analyze(model.A, PoliticalDemocracy)

path.B <- matrix(0, 3, 3)
path.B[1:2, 3] <- NA
path.B[1, 2] <- NA
model.B <- estmodel(LY=loading, BE=path.B, modelType="SEM", indLab=c(paste("x", 1:3, sep=""), paste("y", 1:3, sep="")))

out.B <- analyze(model.B, PoliticalDemocracy)

loading.mis <- matrix("runif(1, -0.2, 0.2)", 11, 3)
loading.mis[is.na(loading)] <- 0

datamodel.A <- model.lavaan(out.A, std=TRUE, LY=loading.mis)
datamodel.B <- model.lavaan(out.B, std=TRUE, LY=loading.mis)

n <- nrow(PoliticalDemocracy)

output.A.A <- sim(20, n=n, model.A, generate=datamodel.A)
output.A.B <- sim(20, n=n, model.B, generate=datamodel.A)
output.B.A <- sim(20, n=n, model.A, generate=datamodel.B)
output.B.B <- sim(20, n=n, model.B, generate=datamodel.B)

# The output may contain some warnings here. When the number of replications increases (e.g., 1000), the warnings
likRatioFit(out.A, out.B, output.A.A, output.A.B, output.B.A, output.B.B)

## End(Not run)
```

| | |
|------------------|--|
| loadingFromAlpha | <i>Find standardized factor loading from coefficient alpha</i> |
|------------------|--|

Description

Find standardized factor loading from coefficient alpha assuming that all items have equal loadings.

Usage

```
loadingFromAlpha(alpha, ni)
```

Arguments

- alpha A desired coefficient alpha value.
- ni A desired number of items.

Value

- result The standardized factor loadings that make desired coefficient alpha with specified number of items.

Author(s)

Sunthud Pornprasertmanit (University of Kansas; <psunthud@ku.edu>)

Examples

```
loadingFromAlpha(0.8, 4)
```

| | |
|------|------------|
| miss | <i>TBA</i> |
|------|------------|

Description

TBA

Usage

```
miss(cov = 0, pmMCAR = 0, pmMAR = 0, nforms = 0, itemGroups = list(),
timePoints = 1, twoMethod = 0, prAttr = 0, package="default", ignoreCols = 0,
threshold = 0, covAsAux = TRUE, logical = NULL, ...)
```


Arguments

| | |
|------------|---|
| cov | Column indices of any normally distributed covariates used in the data set. |
| pmMCAR | Decimal percent of missingness to introduce completely at random on all variables. |
| pmMAR | Decimal percent of missingness to introduce using the listed covariates as predictors. |
| nforms | The number of forms for planned missing data designs, not including the shared form. |
| itemGroups | List of lists of item groupings for planned missing data forms. Without this, items will be divided into groups sequentially (e.g. 1-3,4-6,7-9,10-12) |
| timePoints | Number of timepoints items were measured over. For longitudinal data, planned missing designs will be implemented within each timepoint. |
| twoMethod | With missing on one variable: vector of (column index, percent missing). Will put a given percent missing on that column in the matrix to simulate a two method planned missing data research design. With missing on two or more variables: list of (column indices, percent missing). |
| prAttr | Probability (or vector of probabilities) of an entire case being removed due to attrition at a given time point. See imposeMissing for further details. |
| package | TBA |
| ignoreCols | The columns not imposed any missing values for any missing data patterns |
| threshold | The threshold of covariates that divide between the area to impose missing and the area not to impose missing. The default threshold is the mean of the covariate. |
| covAsAux | If TRUE, the covariate listed in the object will be used as auxiliary variables when putting in the model object. If FALSE, the covariate will be included in the analysis. |
| logical | A matrix of logical values (TRUE/FALSE). If a value in the dataset is corresponding to the TRUE in the logical matrix, the value will be missing. |
| ... | TBA |

Details

TBA

Value

TBA

Author(s)

Alexander M. Schoemann (University of Kansas; <schoemann@ku.edu>), Patrick Miller (University of Notre Dame; <pmille13@nd.edu>), Sunthud Pornprasertmanit (University of Kansas; <psunthud@ku.edu>)

See Also

- [SimMissing](#) for the alternative way to save missing data feature for using in the [sim](#) function.

Examples

```
#Example of imposing 10% MCAR missing in all variables with no imputations (FIML method)
Missing <- miss(pmMCAR=0.1)
summary(Missing)

loading <- matrix(0, 6, 1)
loading[1:6, 1] <- NA
LX <- bind(loading, 0.7)
RPH <- binds(diag(1))
RTD <- binds(diag(6))
CFA.Model <- model(LY = LX, RPS = RPH, RTE = RTD, modelType="CFA")

#Create data
dat <- generate(CFA.Model, n = 20)

#Impose missing
#dat <- run(Missing, dat)

#Analyze data
#out <- run(SimModel, dat)
#summary(out)

#Example to create simMissing object for 3 forms design at 3 timepoints with 10 imputations
Missing <- miss(nforms=3, timePoints=3, numImps=10)
```

model

Data generation template and analysis template for simulation.

Description

Creates a data generation and analysis template (lavaan parameter table) for simulations with structural equation models based on Y-side LISREL design matrices. Each corresponds to a LISREL matrix, but must be a [SimMatrix](#) or [SimVector](#) built using `bind`. In addition to the usual Y-side matrices in LISREL, both PS and TE can be specified using correlations (RPS, RTE) and scaled by a vector of residual variances (VTE, VPS) or total variances (VY, VE). Multiple groups are supported by passing lists of [SimMatrix](#) or [SimVector](#) to arguments, or by specifying the number of groups.

Usage

```
model(LY = NULL, PS = NULL, RPS = NULL, TE = NULL, RTE = NULL, BE = NULL,
VTE = NULL, VY = NULL, VPS = NULL, VE = NULL, TY = NULL, AL = NULL, MY = NULL,
ME = NULL, modelType, indLab=NULL, facLab=NULL, groupLab="group", ngroups=1, smartStart=TRUE)
model.cfa(LY = NULL, PS = NULL, RPS = NULL, TE = NULL, RTE = NULL, VTE = NULL,
VY = NULL, VPS = NULL, VE=NULL, TY = NULL, AL = NULL, MY = NULL, ME = NULL,
indLab=NULL, facLab=NULL, groupLab="group", ngroups=1, smartStart=TRUE)
model.path(PS = NULL, RPS = NULL, BE = NULL, VPS = NULL, VE=NULL, AL = NULL,
ME = NULL, indLab=NULL, facLab=NULL, groupLab="group", ngroups=1, smartStart=TRUE)
model.sem(LY = NULL, PS = NULL, RPS = NULL, TE = NULL, RTE = NULL, BE = NULL,
VTE = NULL, VY = NULL, VPS = NULL, VE=NULL, TY = NULL, AL = NULL, MY = NULL,
ME = NULL, indLab=NULL, facLab=NULL, groupLab="group", ngroups=1, smartStart=TRUE)
```

Arguments

| | |
|------------|---|
| LY | Factor loading matrix from endogenous factors to Y indicators (need to be SimMatrix object). |
| PS | Residual covariance matrix among endogenous factors (need to be SimMatrix object). |
| RPS | Residual correlation matrix among endogenous factors (need to be SimMatrix object). |
| TE | Measurement error covariance matrix among Y indicators (need to be SimMatrix object). |
| RTE | Measurement error correlation matrix among Y indicators (need to be SimMatrix object). |
| BE | Regression coefficient matrix among endogenous factors (need to be SimMatrix object). |
| VTE | Measurement error variance of indicators (need to be SimVector object). |
| VY | Total variance of indicators (need to be SimVector object). NOTE: Either measurement error variance or indicator variance is specified. Both cannot be simultaneously specified. |
| VPS | Residual variance of factors (need to be SimVector object). |
| VE | Total variance of of factors (need to be SimVector object). NOTE: Either residual variance of factors or total variance of factors is specified. Both cannot be simultaneously specified. |
| TY | Measurement intercepts of Y indicators. (need to be SimVector object). |
| AL | Endogenous factor intercept (need to be SimVector object). |
| MY | Overall Y indicator means. (need to be SimVector object). NOTE: Either measurement intercept of indicator mean can be specified. Both cannot be specified simultaneously. |
| ME | Total mean of endogenous factors (need to be SimVector object). NOTE: Either endogenous factor intercept or total mean of endogenous factor is specified. Both cannot be simultaneously specified. |
| modelType | "CFA", "Sem", or "Path". This is specified to ensure that the analysis and data generation template created based on specified matrices in model correspond to what the user intends. |
| indLab | Character vector of indicator labels. If left blank, automatic labels will be generated as y1, y2, ... yy. |
| facLab | Character vector of factor labels. If left blank, automatic labels will be generated as f1, f2, ... ff |
| groupLab | Character of group-variable label (not the names of each group). If left blank, automatic labels will be generated as group |
| ngroups | Integer. Number of groups for data generation, defaults to 1. If larger than one, all specified matrices will be repeated for each additional group. If any matrix argument is a list, the length of this list will be the number of groups and ngroups is ignored. |
| smartStart | Defaults to FALSE. If TRUE, population parameter values that are real numbers will be used as starting values. |

Details

The *simsem* package is intricately tied to the *lavaan* package for analysis of structural equation models. The analysis template that is generated by `model` is a *lavaan* parameter table, a low-level access point to *lavaan* that allows repeated analyses to happen more rapidly. If desired, the parameter table generated can be used directly with *lavaan* for many analyses.

The data generation template is simply a list of *SimMatrix* or *SimVector*. The *SimSem* object can be passed to the function `generate` to generate data.

If multiple group data is desired, the user can optionally either specify the number of groups argument, or pass a list of *SimMatrix* or *SimVector* to any of the matrix arguments. The length of this list will be the number of groups. If only one argument is a list, all other arguments will be automatically replicated to that length, parameters will be identified in the same way, have the same population parameter value/distribution, and have the same misspecification. If only `ngroups` is specified, all arguments will be replicated in this fashion. If equality constraints are present during the automatic replication, these parameters will be constrained to be equal across groups.

The `model.cfa`, `model.path`, and `model.sem` are the shortcuts for the `model` function when `modelType` are "CFA", "Path", and "SEM", respectively.

Value

SimSem object that contains the data generation template (`@dgen`) and analysis template (`@pt`).

Author(s)

Patrick Miller (Univeristy of Notre Dame; <pmille13@end.edu>), Sunthud Pornprasertmanit (University of Kansas; <psunthud@ku.edu>)

See Also

- [sim](#) for simulations using the *SimSem* template.
- [generate](#) To generate data using the *SimSem* template.
- [analyze](#) To analyze real or generated data using the *SimSem* template.
- [draw](#) To draw parameters using the *SimSem* template.

Examples

```
loading <- matrix(0, 6, 2)
loading[1:3, 1] <- NA
loading[4:6, 2] <- NA
LY <- bind(loading, 0.7)

latent.cor <- matrix(NA, 2, 2)
diag(latent.cor) <- 1
RPS <- binds(latent.cor, 0.5)

RTE <- binds(diag(6))

VY <- bind(rep(NA,6),2)

CFA.Model <- model(LY = LY, RPS = RPS, RTE = RTE, modelType = "CFA")
```

| | |
|--------------|--|
| model.lavaan | <i>Build the data generation template and analysis template from the lavaan result</i> |
|--------------|--|

Description

Creates a data generation and analysis template (lavaan parameter table) for simulations with the [lavaan](#) result. Model misspecification may be added into the template by a vector, a matrix, or a list of vectors or matrices (for multiple groups).

Usage

```
model.lavaan(object, std = FALSE, LY = NULL, PS = NULL, RPS = NULL,
  TE = NULL, RTE = NULL, BE = NULL, VTE = NULL, VY = NULL, VPS = NULL,
  VE=NULL, TY = NULL, AL = NULL, MY = NULL, ME = NULL, smartStart=TRUE)
```

Arguments

| | |
|--------|--|
| object | A lavaan object to be used to build the data generation and analysis template. |
| std | If TRUE, use the resulting standardized parameters for data generation. If FALSE, use the unstandardized parameters for data generation. |
| LY | Model misspecification in factor loading matrix from endogenous factors to Y indicators (need to be a matrix or a list of matrices). |
| PS | Model misspecification in residual covariance matrix among endogenous factors (need to be a symmetric matrix or a list of symmetric matrices). |
| RPS | Model misspecification in residual correlation matrix among endogenous factors (need to be a symmetric matrix or a list of symmetric matrices). |
| TE | Model misspecification in measurement error covariance matrix among Y indicators (need to be a symmetric matrix or a list of symmetric matrices). |
| RTE | Model misspecification in measurement error correlation matrix among Y indicators (need to be a symmetric matrix or a list of symmetric matrices). |
| BE | Model misspecification in regression coefficient matrix among endogenous factors (need to be a symmetric matrix or a list of symmetric matrices). |
| VTE | Model misspecification in measurement error variance of indicators (need to be a vector or a list of vectors). |
| VY | Model misspecification in total variance of indicators (need to be a vector or a list of vectors). NOTE: Either measurement error variance or indicator variance is specified. Both cannot be simultaneously specified. |
| VPS | Model misspecification in residual variance of factors (need to be a vector or a list of vectors). |
| VE | Model misspecification in total variance of of factors (need to be a vector or a list of vectors). NOTE: Either residual variance of factors or total variance of factors is specified. Both cannot be simulatneously specified. |
| TY | Model misspecification in measurement intercepts of Y indicators. (need to be a vector or a list of vectors). |
| AL | Model misspecification in endogenous factor intercept (need to be a vector or a list of vectors). |

| | |
|------------|---|
| MY | Model misspecification in overall Y indicator means. (need to be a vector or a list of vectors). NOTE: Either measurement intercept of indicator mean can be specified. Both cannot be specified simultaneously. |
| ME | Model misspecification in total mean of endogenous factors (need to be a vector or a list of vectors). NOTE: Either endogenous factor intercept or total mean of endogenous factor is specified. Both cannot be simultaneously specified. |
| smartStart | Defaults to FALSE. If TRUE, population parameter values that are real numbers will be used as starting values. |

Value

SimSem object that contains the data generation template (@dgen) and analysis template (@pt).

Author(s)

Sunthud Pornprasertmanit (University of Kansas; <psunthud@ku.edu>)

See Also

- [model](#) To build data generation and data analysis template for simulation.
- [sim](#) for simulations using the [SimSem](#) template.
- [generate](#) To generate data using the [SimSem](#) template.
- [analyze](#) To analyze real or generated data using the [SimSem](#) template.
- [draw](#) To draw parameters using the [SimSem](#) template.

Examples

```
HS.model <- ' visual =~ x1 + x2 + x3
              textual =~ x4 + x5 + x6
              speed  =~ x7 + x8 + x9 '

fit <- cfa(HS.model, data=HolzingerSwineford1939)
datamodel <- model.lavaan(fit, std=TRUE)
```

| | |
|------------------|---|
| multipleAllEqual | <i>Test whether all objects are equal</i> |
|------------------|---|

Description

Test whether all objects are equal. The test is based on the [all.equal](#) function.

Usage

```
multipleAllEqual(...)
```

Arguments

... The target objects

Value

TRUE if all objects are equal.

Author(s)

Sunthud Pornprasertmanit (University of Kansas; <psunthud@ku.edu>)

Examples

```
multipleAllEqual(1:5, 1:5, seq(2, 10, 2)/2)
multipleAllEqual(1:5, 1:6, seq(2, 10, 2)/2)
```

| | |
|-------------|------------------------------------|
| overlapHist | <i>Plot overlapping histograms</i> |
|-------------|------------------------------------|

Description

Plot overlapping histograms

Usage

```
overlapHist(a, b, colors=c("red","blue","purple"), breaks=NULL, xlim=NULL,
ylim=NULL, main=NULL, xlab=NULL, swap=FALSE)
```

Arguments

| | |
|--------|---|
| a | Data for the first histogram |
| b | Data for the second histogram |
| colors | Colors for the first histogram, the second histogram, and the overlapping areas. |
| breaks | How many breaks users used in each histogram (should not be used) |
| xlim | The range of x-axis |
| ylim | The range of y-axis |
| main | The title of the figure |
| xlab | The labels of x-axis |
| swap | Specify TRUE to plot b first and then a. The default is FALSE to plot a first and then b. |

Value

None. This function will plot only.

Author(s)

Chris Miller provided this code on <http://chrisamiller.com/science/2010/07/20/transparent-overlapping-h>
The code is modified by Sunthud Pornprasertmanit (University of Kansas; psunthud@ku.edu)

Examples

```
# This function is not a public function.

# a <- rnorm(10000, 0, 1)
# b <- rnorm(10000, 1, 1.5)
# overlapHist(a, b, main="Example")
```

| | |
|-------------|---|
| plot3DQtile | <i>Build a persepctive plot or contour plot of a quantile of predicted values</i> |
|-------------|---|

Description

Build a persepctive plot or contour plot of a quantile of predicted values

Usage

```
plot3DQtile(x, y, z, df=0, qtile=0.5, useContour=TRUE, xlab=NULL,
ylab=NULL, zlab=NULL, main=NULL)
```

Arguments

| | |
|------------|---|
| x | The values of the first variable (e.g., a vector of sample size) |
| y | The values of the second variable (e.g., a vector of percent missing) |
| z | The values of the dependent variable |
| df | The degree of freedom in spline method |
| qtile | The quantile values used to plot a graph |
| useContour | If TRUE, use contour plot. If FALSE, use perspective plot. |
| xlab | The labels of x-axis |
| ylab | The labels of y-axis |
| zlab | The labels of z-axis |
| main | The title of the graph |

Value

None. This function will plot only.

Author(s)

Sunthud Pornprasertmanit (University of Kansas; psunthud@ku.edu)

Examples

```
# No example
```


plotCutoff

*Plot sampling distributions of fit indices with fit indices cutoffs***Description**

This function will plot sampling distributions of null hypothesis fit indices. The users may add cutoffs by specifying the alpha level.

Usage

```
plotCutoff(object, ...)
```

Arguments

| | |
|--------|---|
| object | The object (SimResult or <code>data.frame</code>) that contains values of fit indices in each distribution. |
| ... | Other arguments specific to different types of object you pass in the function. |

Value

NONE. Only plot the fit indices distributions.

Details in ...

- `cutoff`: A priori cutoffs for fit indices, saved in a vector
- `cutoff2`: Another set of priori cutoffs for fit indices, saved in a vector
- `alpha`: A priori alpha level to getCutoffs of fit indices (do not specify when you have cutoff)
- `revDirec`: The default is to find critical point on the side that indicates worse fit (the right side of RMSEA or the left side of CFI). If specifying as TRUE, the directions are reversed.
- `usedFit`: The name of fit indices that researchers wish to plot
- `useContour`: If there are two of sample size, percent completely at random, and percent missing at random are varying, the `plotCutoff` function will provide 3D graph. Contour graph is a default. However, if this is specified as FALSE, perspective plot is used.

Author(s)

Sunthud Pornprasertmanit (University of Kansas; <psunthud@ku.edu>)

See Also

- [SimResult](#) for `simResult` that used in this function.
- [getCutoff](#) to find values of cutoffs based on null hypothesis sampling distributions only

Examples

```
## Not run:
loading <- matrix(0, 6, 2)
loading[1:3, 1] <- NA
loading[4:6, 2] <- NA
loadingValues <- matrix(0, 6, 2)
loadingValues[1:3, 1] <- 0.7
loadingValues[4:6, 2] <- 0.7
LX <- bind(loading, loadingValues)
latent.cor <- matrix(NA, 2, 2)
diag(latent.cor) <- 1
RPH <- binds(latent.cor, 0.5)
error.cor <- matrix(0, 6, 6)
diag(error.cor) <- 1
RTD <- binds(error.cor)
CFA.Model <- model(LY = LX, RPS = RPH, RTE = RTD, modelType="CFA")
# We make the examples running only 5 replications to save time.
# In reality, more replications are needed.
Output <- sim(5, n=200, model=CFA.Model)
plotCutoff(Output, 0.05, usedFit=c("RMSEA", "SRMR", "CFI", "TLI"))

# Varying N
Output2 <- sim(NULL, n=seq(450, 500, 10), model=CFA.Model)
plotCutoff(Output2, 0.05)

# Varying N and pmMCAR
Output3 <- sim(NULL, n=seq(450, 500, 10), pmMCAR=c(0, 0.05, 0.1, 0.15), model=CFA.Model)
plotCutoff(Output3, 0.05)

## End(Not run)
```

| | |
|------------------|---|
| plotCutoffNested | <i>Plot sampling distributions of the differences in fit indices between nested models with fit indices cutoffs</i> |
|------------------|---|

Description

This function will plot sampling distributions of the differences in fit indices between nested models if the nested model is true. The users may add cutoffs by specifying the alpha level.

Usage

```
plotCutoffNested(nested, parent, alpha = 0.05, cutoff = NULL,
  usedFit = NULL, useContour = T)
```

Arguments

| | |
|--------|--|
| nested | SimResult that saves the analysis results of nested model from multiple replications |
| parent | SimResult that saves the analysis results of parent model from multiple replications |
| alpha | A priori alpha level |

| | |
|------------|--|
| cutoff | A priori cutoffs for fit indices, saved in a vector |
| usedFit | Vector of names of fit indices that researchers wish to plot the sampling distribution. |
| useContour | If there are two of sample size, percent completely at random, and percent missing at random are varying, the plotCutoff function will provide 3D graph. Contour graph is a default. However, if this is specified as FALSE, perspective plot is used. |

Value

NONE. Only plot the fit indices distributions.

Author(s)

Sunthud Pornprasertmanit (University of Kansas; <psunthud@ku.edu>)

See Also

- [SimResult](#) for simResult that used in this function.
- [getCutoffNested](#) to find the difference in fit indices cutoffs

Examples

```
## Not run:
loading.null <- matrix(0, 6, 1)
loading.null[1:6, 1] <- NA
LX.NULL <- bind(loading.null, 0.7)
RPH.NULL <- binds(diag(1))
RTD <- binds(diag(6))
CFA.Model.NULL <- model(LY = LX.NULL, RPS = RPH.NULL, RTE = RTD, modelType="CFA")

loading.alt <- matrix(0, 6, 2)
loading.alt[1:3, 1] <- NA
loading.alt[4:6, 2] <- NA
LX.ALT <- bind(loading.alt, 0.7)
latent.cor.alt <- matrix(NA, 2, 2)
diag(latent.cor.alt) <- 1
RPH.ALT <- binds(latent.cor.alt, "runif(1, 0.7, 0.9)")
CFA.Model.ALT <- model(LY = LX.ALT, RPS = RPH.ALT, RTE = RTD, modelType="CFA")

# The actual number of replications should be greater than 10.
Output.NULL.NULL <- sim(10, n=500, model=CFA.Model.NULL)
Output.NULL.ALT <- sim(10, n=500, model=CFA.Model.ALT, generate=CFA.Model.NULL)

plotCutoffNested(Output.NULL.NULL, Output.NULL.ALT, alpha=0.05)

## End(Not run)
```

| | |
|---------------------|---|
| plotCutoffNonNested | <i>Plot sampling distributions of the differences in fit indices between non-nested models with fit indices cutoffs</i> |
|---------------------|---|

Description

This function will plot sampling distributions of the differences in fit indices between non-nested models. The users may add cutoffs by specifying the alpha level.

Usage

```
plotCutoffNonNested(dat1Mod1, dat1Mod2, dat2Mod1=NULL, dat2Mod2=NULL,
alpha=0.05, cutoff = NULL, usedFit = NULL, useContour = T, onetailed=FALSE)
```

Arguments

| | |
|------------|--|
| dat1Mod1 | SimResult that saves the simulation of analyzing Model 1 by datasets created from Model 1 |
| dat1Mod2 | SimResult that saves the simulation of analyzing Model 2 by datasets created from Model 1 |
| dat2Mod1 | SimResult that saves the simulation of analyzing Model 1 by datasets created from Model 2 |
| dat2Mod2 | SimResult that saves the simulation of analyzing Model 2 by datasets created from Model 2 |
| alpha | A priori alpha level |
| cutoff | A priori cutoffs for fit indices, saved in a vector |
| usedFit | Vector of names of fit indices that researchers wish to plot the sampling distribution. |
| useContour | If there are two of sample size, percent completely at random, and percent missing at random are varying, the plotCutoff function will provide 3D graph. Contour graph is a default. However, if this is specified as FALSE, perspective plot is used. |
| onetailed | If TRUE, the function will find the cutoff from one-tail test. If FALSE, the function will find the cutoff from two-tailed test. |

Value

NONE. Only plot the fit indices distributions.

Author(s)

Sunthud Pornprasertmanit (University of Kansas; <psunthud@ku.edu>)

See Also

- [SimResult](#) for simResult that used in this function.
- [getCutoffNonNested](#) to find the difference in fit indices cutoffs for non-nested model comparison

Examples

```
## Not run:
loading.A <- matrix(0, 8, 2)
loading.A[1:3, 1] <- NA
loading.A[4:8, 2] <- NA
LX.A <- bind(loading.A, 0.7)
latent.cor <- matrix(NA, 2, 2)
diag(latent.cor) <- 1
RPH <- binds(latent.cor, "runif(1, 0.7, 0.9)")
RTD <- binds(diag(8))
CFA.Model.A <- model(LY = LX.A, RPS = RPH, RTE = RTD, modelType="CFA")

loading.B <- matrix(0, 8, 2)
loading.B[1:4, 1] <- NA
loading.B[5:8, 2] <- NA
LX.B <- bind(loading.B, 0.7)
CFA.Model.B <- model(LY = LX.B, RPS = RPH, RTE = RTD, modelType="CFA")

# The actual number of replications should be greater than 10.
Output.A.A <- sim(10, n=500, model=CFA.Model.A, generate=CFA.Model.A)
Output.A.B <- sim(10, n=500, model=CFA.Model.B, generate=CFA.Model.A)
Output.B.A <- sim(10, n=500, model=CFA.Model.A, generate=CFA.Model.B)
Output.B.B <- sim(10, n=500, model=CFA.Model.B, generate=CFA.Model.B)

plotCutoffNonNested(Output.A.A, Output.A.B, Output.B.A, Output.B.B)
plotCutoffNonNested(Output.A.A, Output.A.B)
plotCutoffNonNested(Output.A.A, Output.A.B, onetailed=TRUE)

## End(Not run)
```

plotDist

Plot a distribution of a data distribution object

Description

Plot a distribution of a data distribution object

Usage

```
plotDist(object, xlim = NULL, ylim = NULL, r = 0, var = NULL, contour = TRUE)
```

Arguments

| | |
|--------|--|
| object | The data distribution object (SimDataDist) to plot a distribution |
| xlim | A numeric vector with two elements specifying the lower and upper limit of the x-axis to be plotted. |
| ylim | A numeric vector with two elements specifying the lower and upper limit of the y-axis to be plotted. This argument is applicable for the joint distribution of two dimensions only |
| r | The correlation of two dimensions in the joint distribution |
| var | A vector of the index of variables to be plotted. The length of vector cannot be greater than 2. |

contour Applicable if two variables are used only. If TRUE, the contour plot is provided. If FALSE, the perspective plot is provided.

Value

No return value. This function will plot a graph only.

Author(s)

Sunthud Pornprasertmanit (University of Kansas; <psunthud@ku.edu>)

See Also

- [SimDataDist](#) for plotting a data distribution object

Examples

```
datadist <- bindDist(c("chisq", "t", "f"), list(df=5), list(df=3), list(df1=3, df2=5))
plotDist(datadist, r=0.5, var=1:2)
plotDist(datadist, var=3)
```

plotIndividualScatter *Plot an overlaying scatter plot visualizing the power of rejecting misspecified models*

Description

Plot the fit indices value against the value of predictors. The plot will include the fit indices value of the alternative models, the fit indices value of the null model (if specified), and the fit indices cutoffs (if specified).

Usage

```
plotIndividualScatter(altVec, nullVec=NULL, cutoff=NULL, x, main = NULL)
```

Arguments

| | |
|---------|---|
| altVec | The vector saving the fit index distribution when the hypothesized model is FALSE. |
| nullVec | The vector saving the fit index distribution when the hypothesized model is TRUE. |
| cutoff | A priori cutoff |
| x | The data.frame of the predictor values. The number of rows of the x argument should be equal to the number of rows in the object. |
| main | The title of the graph |

Value

NONE. Only plot the fit indices distributions.

Author(s)

Sunthud Pornprasertmanit (University of Kansas; <psunthud@ku.edu>)

See Also

- [plotPowerFit](#)

Examples

```
# No example
```

| | |
|-----------------|---|
| plotLogisticFit | <i>Plot multiple logistic curves for predicting whether rejecting a mis-specified model</i> |
|-----------------|---|

Description

This function will find the fit indices cutoff values if not specified, then check whether the hypothesized model is rejected in each dataset, and plot the logistic curve given the value of predictors.

Usage

```
plotLogisticFit(altObject, nullObject=NULL, cutoff=NULL,
  usedFit=NULL, x, xval, alpha=0.05, useContour=TRUE, df=0)
```

Arguments

| | |
|------------|--|
| altObject | The result object (data.frame) saves the simulation result of fitting the hypothesized model when the hypothesized model is FALSE. |
| nullObject | The result object (data.frame) saves the simulation result of fitting the hypothesized model when the hypothesized model is TRUE. This argument may be not specified if the cutoff is specified. |
| cutoff | A vector of priori cutoffs for fit indices. |
| usedFit | Vector of names of fit indices that researchers wish to plot. |
| alpha | A priori alpha level |
| x | The data.frame of the predictor values. The number of rows of the x argument should be equal to the number of rows in the object. |
| xval | The values of predictor that researchers would like to find the fit indices cutoffs from. |
| useContour | If there are two of sample size, percent completely at random, and percent missing at random are varying, the plotCutoff function will provide 3D graph. Contour graph is a default. However, if this is specified as FALSE, perspective plot is used. |
| df | The degree of freedom used in spline method in predicting the fit indices by the predictors. If df is 0, the spline method will not be applied. |

Value

NONE. Only plot the fit indices distributions.

Author(s)

Sunthud Pornprasertmanit (University of Kansas; <psunthud@ku.edu>)

See Also

- [plotPowerFit](#)

Examples

```
# No example
```

| | |
|------------|--|
| plotMisfit | <i>Plot the population misfit in the result object</i> |
|------------|--|

Description

Plot a histogram of the amount of population misfit in parameter result object or the scatter plot of the relationship between misspecified parameter and the population misfit or the fit indices

Usage

```
plotMisfit(object, usedFit="default", misParam=NULL)
```

Arguments

| | |
|----------|---|
| object | The result object, SimResult |
| usedFit | The sample fit indices or population misfit used to plot. All sample fit indices are available. The available population misfit are "pop.f0", "pop.rmsea", and "pop.srmr". If the misParam is not specified, all population misfit are used. If the misParam is specified, the "pop.rmsea" is used in the plot. |
| misParam | The index or the name of misspecified parameters used to plot. |

Value

None. This function will plot only.

Author(s)

Sunthud Pornprasertmanit (University of Kansas; <psunthud@ku.edu>)

Examples

```
path.BE <- matrix(0, 4, 4)
path.BE[3, 1:2] <- NA
path.BE[4, 3] <- NA
starting.BE <- matrix("", 4, 4)
starting.BE[3, 1:2] <- "runif(1, 0.3, 0.5)"
starting.BE[4, 3] <- "runif(1, 0.5, 0.7)"
mis.path.BE <- matrix(0, 4, 4)
mis.path.BE[4, 1:2] <- "runif(1, -0.1, 0.1)"
BE <- bind(path.BE, starting.BE, misspec=mis.path.BE)
```



```

residual.error <- diag(4)
residual.error[1,2] <- residual.error[2,1] <- NA
RPS <- binds(residual.error, "rnorm(1, 0.3, 0.1)")

ME <- bind(rep(NA, 4), 0)

Path.Model <- model(RPS = RPS, BE = BE, ME = ME, modelType="Path")

# The number of replications in actual analysis should be much more than 5
ParamObject <- sim(20, n=500, Path.Model)
plotMisfit(ParamObject)

plotMisfit(ParamObject, misParam=1:2)

```

| | |
|--------------|---|
| plotOverHist | <i>Plot multiple overlapping histograms</i> |
|--------------|---|

Description

Plot multiple overlapping histograms and find the cutoff values if not specified

Usage

```
plotOverHist(altObject, nullObject, cutoff=NULL, usedFit=NULL, alpha=0.05,
cutoff2=NULL, cutoff3=NULL, cutoff4=NULL)
```

Arguments

| | |
|------------|--|
| altObject | The result object (data.frame) saves the simulation result of fitting the hypothesized model when the hypothesized model is FALSE. |
| nullObject | The result object (data.frame) saves the simulation result of fitting the hypothesized model when the hypothesized model is TRUE. |
| cutoff | A vector of priori cutoffs for fit indices. |
| usedFit | Vector of names of fit indices that researchers wish to plot. |
| alpha | A priori alpha level |
| cutoff2 | Another vector of priori cutoffs for fit indices. |
| cutoff3 | A vector of priori cutoffs for fit indices for the altObject. |
| cutoff4 | Another vector of priori cutoffs for fit indices for the altObject. |

Value

NONE. Only plot the fit indices distributions.

Author(s)

Sunthud Pornprasertmanit (University of Kansas; <psunthud@ku.edu>)

See Also

- [plotPowerFit](#)

Examples

```
# No example
```

plotPower

Make a power plot of a parameter given varying parameters

Description

Make a power plot of a parameter given varying parameters (e.g., sample size, percent missing completely at random, or random parameters in the model)

Usage

```
plotPower(object, powerParam, alpha = 0.05, contParam = NULL, contN = TRUE,
  contMCAR = TRUE, contMAR = TRUE, useContour=TRUE)
```

Arguments

| | |
|------------|---|
| object | SimResult that includes at least one randomly varying parameter (e.g. sample size, percent missing, model parameters) |
| powerParam | Vector of parameters names that the user wishes to find power for. This can be a vector of names (e.g., "LY1_1", "LY2_2"). |
| alpha | Alpha level to use for power analysis. |
| contParam | Vector of parameters names that vary over replications that users wish to use in the plot. |
| contN | Include the varying sample size in the power plot if available |
| contMCAR | Include the varying MCAR (missing completely at random percentage) in the power plot if available |
| contMAR | Include the varying MAR (missing at random percentage) in the power plot if available |
| useContour | This argument is used when users specify to plot two varying parameters. If TRUE, the contour plot is used. If FALSE, perspective plot is used. |

Details

Predicting whether each replication is significant or not by varying parameters using logistic regression (without interaction). Then, plot the logistic curves predicting the probability of significance against the target varying parameters.

Value

Not return any value. This function will plot a graph only.

Author(s)

Sunthud Pornprasertmanit (University of Kansas; <psunthud@ku.edu>), Alexander M. Schoemann (University of Kansas; <schoemann@ku.edu>)

See Also

- [SimResult](#) to see how to create a simResult object with randomly varying parameters.
- [getPower](#) to obtain a statistical power given varying parameters values.

Examples

```
## Not run:
# Specify Sample Size by n
loading <- matrix(0, 6, 1)
loading[1:6, 1] <- NA
LX <- bind(loading, 0.4)
RPH <- binds(diag(1))
RTD <- binds(diag(6))
CFA.Model <- model(LY = LX, RPS = RPH, RTE = RTD, modelType="CFA")

# Specify both sample size and percent missing completely at random
Output <- sim(NULL, n=seq(100, 200, 20), pmMCAR=c(0, 0.1, 0.2), model=CFA.Model)
plotPower(Output, "1.LY1_1", contMCAR=FALSE)

## End(Not run)
```

plotPowerFit

Plot sampling distributions of fit indices that visualize power of rejecting datasets underlying misspecified models

Description

This function will plot sampling distributions of fit indices that visualize power in rejecting the misspecified models

Usage

```
plotPowerFit(altObject, nullObject = NULL, cutoff = NULL, usedFit = NULL,
alpha = 0.05, contN = TRUE, contMCAR = TRUE, contMAR = TRUE,
useContour = TRUE, logistic = TRUE)
```

Arguments

| | |
|------------|---|
| altObject | The result object (SimResult) saves the simulation result of fitting the hypothesized model when the hypothesized model is FALSE. |
| nullObject | The result object (SimResult) saves the simulation result of fitting the hypothesized model when the hypothesized model is TRUE. This argument may be not specified if the cutoff is specified. |
| cutoff | A vector of priori cutoffs for fit indices. |
| usedFit | Vector of names of fit indices that researchers wish to plot. |
| alpha | A priori alpha level |
| contN | Include the varying sample size in the power plot if available |
| contMCAR | Include the varying MCAR (missing completely at random percentage) in the power plot if available |

| | |
|------------|---|
| contMAR | Include the varying MAR (missing at random percentage) in the power plot if available |
| useContour | If there are two of sample size, percent completely at random, and percent missing at random are varying, the plotCutoff function will provide 3D graph. Contour graph is a default. However, if this is specified as FALSE, perspective plot is used. |
| logistic | If logistic is TRUE and the varying parameter exists (e.g., sample size or percent missing), the plot based on logistic regression predicting the significance by the varying parameters is preferred. If FALSE, the overlaying scatterplot with a line of cutoff is plotted. |

Value

NONE. Only plot the fit indices distributions.

Author(s)

Sunthud Pornprasertmanit (University of Kansas; <psunthud@ku.edu>)

See Also

- [SimResult](#) for simResult that used in this function.
- [getCutoff](#) to find values of cutoffs based on null hypothesis sampling distributions only
- [getPowerFit](#) to find power of rejecting the hypothesized model when the hypothesized model is FALSE.

Examples

```
## Not run:
loading.null <- matrix(0, 6, 1)
loading.null[1:6, 1] <- NA
LX.NULL <- bind(loading.null, 0.7)
RPH.NULL <- binds(diag(1))
RTD <- binds(diag(6))
CFA.Model.NULL <- model(LY = LX.NULL, RPS = RPH.NULL, RTE = RTD, modelType="CFA")
# We make the examples running only 5 replications to save time.
# In reality, more replications are needed.
Output.NULL <- sim(50, n=50, model=CFA.Model.NULL, generate=CFA.Model.NULL)

loading.alt <- matrix(0, 6, 2)
loading.alt[1:3, 1] <- NA
loading.alt[4:6, 2] <- NA
LX.ALT <- bind(loading.alt, 0.7)
latent.cor.alt <- matrix(NA, 2, 2)
diag(latent.cor.alt) <- 1
RPH.ALT <- binds(latent.cor.alt, 0.5)
CFA.Model.ALT <- model(LY = LX.ALT, RPS = RPH.ALT, RTE = RTD, modelType="CFA")
Output.ALT <- sim(50, n=50, model=CFA.Model.NULL, generate=CFA.Model.ALT)

datNull <- generate(CFA.Model.NULL, n=50, params=TRUE)
datAlt <- generate(CFA.Model.ALT, n=50, params=TRUE)
outNull <- analyze(CFA.Model.NULL, datNull)
outAlt <- analyze(CFA.Model.NULL, datAlt)
summaryFit(Output.NULL)
```

```
summaryFit(Output.ALT)

plotPowerFit(Output.ALT, nullObject=Output.NULL, alpha=0.05, usedFit=c("RMSEA", "CFI", "TLI", "SRMR"))
Rule.of.thumb <- c(RMSEA=0.05, CFI=0.95, TLI=0.95, SRMR=0.06)
plotPowerFit(Output.ALT, cutoff=Rule.of.thumb, alpha=0.05, usedFit=c("RMSEA", "CFI", "TLI", "SRMR"))

Output.NULL2 <- sim(NULL, n=seq(50, 250, 25), model=CFA.Model.NULL, generate=CFA.Model.NULL)
Output.ALT2 <- sim(NULL, n=seq(50, 250, 25), model=CFA.Model.NULL, generate=CFA.Model.ALT)

plotPowerFit(Output.ALT2, nullObject=Output.NULL2, alpha=0.05, usedFit=c("RMSEA", "CFI", "TLI", "SRMR"))
plotPowerFit(Output.ALT2, cutoff=Rule.of.thumb, alpha=0.05, usedFit=c("RMSEA", "CFI", "TLI", "SRMR"))

## End(Not run)
```

| | |
|----------------|---|
| plotPowerFitDf | <i>Plot sampling distributions of fit indices that visualize power of rejecting datasets underlying misspecified models</i> |
|----------------|---|

Description

This function will plot sampling distributions of fit indices that visualize power in rejecting the misspecified models. This function is similar to the [plotPowerFit](#) function but the input distributions are data.frame.

Usage

```
plotPowerFitDf(altObject, nullObject = NULL, cutoff = NULL, usedFit = NULL, alpha = 0.05, x = NULL)
```

Arguments

| | |
|------------|---|
| altObject | The result object (data.frame) saves the simulation result of fitting the hypothesized model when the hypothesized model is FALSE. |
| nullObject | The result object (data.frame) saves the simulation result of fitting the hypothesized model when the hypothesized model is TRUE. This argument may be not specified if the cutoff is specified. |
| cutoff | A vector of priori cutoffs for fit indices. |
| usedFit | Vector of names of fit indices that researchers wish to plot. |
| alpha | A priori alpha level |
| x | The data.frame of the predictor values. The number of rows of the x argument should be equal to the number of rows in the object. |
| xval | The values of predictor that researchers would like to find the fit indices cutoffs from. |
| useContour | If there are two of sample size, percent completely at random, and percent missing at random are varying, the plotCutoff function will provide 3D graph. Contour graph is a default. However, if this is specified as FALSE, perspective plot is used. |
| logistic | If logistic is TRUE and the varying parameter exists (e.g., sample size or percent missing), the plot based on logistic regression predicting the significance by the varying parameters is preferred. If FALSE, the overlaying scatterplot with a line of cutoff is plotted. |

Value

NONE. Only plot the fit indices distributions.

Author(s)

Sunthud Pornprasertmanit (University of Kansas; <psunthud@ku.edu>)

See Also

- [plotPowerFit](#)

Examples

```
# No example
```

| | |
|--------------------|--|
| plotPowerFitNested | <i>Plot power of rejecting a nested model in a nested model comparison by each fit index</i> |
|--------------------|--|

Description

This function will plot sampling distributions of the differences in fit indices between parent and nested models. Two sampling distributions will be compared: nested model is FALSE (alternative model) and nested model is TRUE (null model).

Usage

```
plotPowerFitNested(altNested, altParent, nullNested = NULL,
  nullParent = NULL, cutoff = NULL, usedFit = NULL, alpha = 0.05,
  contN = TRUE, contMCAR = TRUE, contMAR = TRUE, useContour = TRUE,
  logistic = TRUE)
```

Arguments

| | |
|------------|--|
| altNested | SimResult that saves the simulation result of the nested model when the nested model is FALSE. |
| altParent | SimResult that saves the simulation result of the parent model when the nested model is FALSE. |
| nullNested | SimResult that saves the simulation result of the nested model when the nested model is TRUE. This argument may not be specified if the cutoff is specified. |
| nullParent | SimResult that saves the simulation result of the parent model when the nested model is TRUE. This argument may not be specified if the cutoff is specified. |
| cutoff | A vector of priori cutoffs for the differences in fit indices. |
| usedFit | Vector of names of fit indices that researchers wish to plot. |
| alpha | A priori alpha level |
| contN | Include the varying sample size in the power plot if available |
| contMCAR | Include the varying MCAR (missing completely at random percentage) in the power plot if available |

| | |
|------------|---|
| contMAR | Include the varying MAR (missing at random percentage) in the power plot if available |
| useContour | If there are two of sample size, percent completely at random, and percent missing at random are varying, the plotCutoff function will provide 3D graph. Contour graph is a default. However, if this is specified as FALSE, perspective plot is used. |
| logistic | If logistic is TRUE and the varying parameter exists (e.g., sample size or percent missing), the plot based on logistic regression predicting the significance by the varying parameters is preferred. If FALSE, the overlaying scatterplot with a line of cutoff is plotted. |

Value

NONE. Only plot the fit indices distributions.

Author(s)

Sunthud Pornprasertmanit (University of Kansas; <psunthud@ku.edu>)

See Also

- [SimResult](#) for simResult that used in this function.
- [getCutoffNested](#) to find the cutoffs of the differences in fit indices
- [plotCutoffNested](#) to visualize the cutoffs of the differences in fit indices
- [getPowerFitNested](#) to find the power in rejecting the nested model by the difference in fit indices cutoffs

Examples

```
## Not run:
loading.null <- matrix(0, 6, 1)
loading.null[1:6, 1] <- NA
LX.NULL <- bind(loading.null, 0.7)
RPH.NULL <- binds(diag(1))
RTD <- binds(diag(6))
CFA.Model.NULL <- model(LY = LX.NULL, RPS = RPH.NULL, RTE = RTD, modelType="CFA")
```

```
loading.alt <- matrix(0, 6, 2)
loading.alt[1:3, 1] <- NA
loading.alt[4:6, 2] <- NA
LX.ALT <- bind(loading.alt, 0.7)
latent.cor.alt <- matrix(NA, 2, 2)
diag(latent.cor.alt) <- 1
RPH.ALT <- binds(latent.cor.alt, 0.7)
CFA.Model.ALT <- model(LY = LX.ALT, RPS = RPH.ALT, RTE = RTD, modelType="CFA")
```

```
Output.NULL.NULL <- sim(10, n=500, model=CFA.Model.NULL, generate=CFA.Model.NULL)
Output.ALT.NULL <- sim(10, n=500, model=CFA.Model.NULL, generate=CFA.Model.ALT)
Output.NULL.ALT <- sim(10, n=500, model=CFA.Model.ALT, generate=CFA.Model.NULL)
Output.ALT.ALT <- sim(10, n=500, model=CFA.Model.ALT, generate=CFA.Model.ALT)
```

```
plotPowerFitNested(Output.ALT.NULL, Output.ALT.ALT, nullNested=Output.NULL.NULL, nullParent=Output.NULL.ALT)
plotPowerFitNested(Output.ALT.NULL, Output.ALT.ALT, nullNested=Output.NULL.NULL, nullParent=Output.NULL.ALT)
```

```

Output.NULL.NULL2 <- sim(NULL, n=seq(50, 500, 25), model=CFA.Model.NULL, generate=CFA.Model.NULL)
Output.ALT.NULL2 <- sim(NULL, n=seq(50, 500, 25), model=CFA.Model.NULL, generate=CFA.Model.ALT)
Output.NULL.ALT2 <- sim(NULL, n=seq(50, 500, 25), model=CFA.Model.ALT, generate=CFA.Model.NULL)
Output.ALT.ALT2 <- sim(NULL, n=seq(50, 500, 25), model=CFA.Model.ALT, generate=CFA.Model.ALT)

plotPowerFitNested(Output.ALT.NULL2, Output.ALT.ALT2, nullNested=Output.NULL.NULL2, nullParent=Output.NULL.NULL2)

plotPowerFitNested(Output.ALT.NULL2, Output.ALT.ALT2, nullNested=Output.NULL.NULL2, nullParent=Output.NULL.NULL2)

plotPowerFitNested(Output.ALT.NULL2, Output.ALT.ALT2, cutoff=c(CFI=-0.1), logistic=FALSE)

## End(Not run)

```

`plotPowerFitNonNested` *Plot power of rejecting a non-nested model based on a difference in fit index*

Description

Plot the proportion of the difference in fit indices from one model that does not in the range of sampling distribution from another model (reject that the dataset comes from the second model) or indicates worse fit than a specified cutoff. This plot can show the proportion in the second model that does not in the range of sampling distribution from the first model too.

Usage

```

plotPowerFitNonNested(dat2Mod1, dat2Mod2, dat1Mod1=NULL, dat1Mod2=NULL,
  cutoff = NULL, usedFit = NULL, alpha = 0.05, contN = TRUE, contMCAR = TRUE,
  contMAR = TRUE, useContour = TRUE, logistic = TRUE, onetailed = FALSE)

```

Arguments

| | |
|-----------------------|---|
| <code>dat2Mod1</code> | SimResult that saves the simulation of analyzing Model 1 by datasets created from Model 2 |
| <code>dat2Mod2</code> | SimResult that saves the simulation of analyzing Model 2 by datasets created from Model 2 |
| <code>dat1Mod1</code> | SimResult that saves the simulation of analyzing Model 1 by datasets created from Model 1 |
| <code>dat1Mod2</code> | SimResult that saves the simulation of analyzing Model 2 by datasets created from Model 1 |
| <code>cutoff</code> | A vector of priori cutoffs for the differences in fit indices. |
| <code>usedFit</code> | Vector of names of fit indices that researchers wish to plot. |
| <code>alpha</code> | A priori alpha level |
| <code>contN</code> | Include the varying sample size in the power plot if available |
| <code>contMCAR</code> | Include the varying MCAR (missing completely at random percentage) in the power plot if available |
| <code>contMAR</code> | Include the varying MAR (missing at random percentage) in the power plot if available |

| | |
|------------|---|
| useContour | If there are two of sample size, percent completely at random, and percent missing at random are varying, the plotCutoff function will provide 3D graph. Contour graph is a default. However, if this is specified as FALSE, perspective plot is used. |
| logistic | If logistic is TRUE and the varying parameter exists (e.g., sample size or percent missing), the plot based on logistic regression predicting the significance by the varying parameters is preferred. If FALSE, the overlaying scatterplot with a line of cutoff is plotted. |
| onetailed | If TRUE, the function will use the cutoff from one-tail test. If FALSE, the function will use the cutoff from two-tailed test. |

Value

NONE. Only plot the fit indices distributions.

Author(s)

Sunthud Pornprasertmanit (University of Kansas; <psunthud@ku.edu>)

See Also

- [SimResult](#) for simResult that used in this function.
- [getCutoffNonNested](#) to find the cutoffs of the differences in fit indices for non-nested model comparison
- [plotCutoffNonNested](#) to visualize the cutoffs of the differences in fit indices for non-nested model comparison
- [getPowerFitNonNested](#) to find the power in rejecting the non-nested model by the difference in fit indices cutoffs

Examples

```
## Not run:
loading.A <- matrix(0, 8, 2)
loading.A[1:3, 1] <- NA
loading.A[4:8, 2] <- NA
LX.A <- bind(loading.A, 0.7)
latent.cor <- matrix(NA, 2, 2)
diag(latent.cor) <- 1
RPH <- binds(latent.cor, "runif(1, 0.7, 0.9)")
RTD <- binds(diag(8))
CFA.Model.A <- model(LY = LX.A, RPS = RPH, RTE = RTD, modelType="CFA")

loading.B <- matrix(0, 8, 2)
loading.B[1:4, 1] <- NA
loading.B[5:8, 2] <- NA
LX.B <- bind(loading.B, 0.7)
CFA.Model.B <- model(LY = LX.B, RPS = RPH, RTE = RTD, modelType="CFA")

# The actual number of replications should be greater than 10.
Output.A.A <- sim(10, n=500, model=CFA.Model.A, generate=CFA.Model.A)
Output.A.B <- sim(10, n=500, model=CFA.Model.B, generate=CFA.Model.A)
Output.B.A <- sim(10, n=500, model=CFA.Model.A, generate=CFA.Model.B)
Output.B.B <- sim(10, n=500, model=CFA.Model.B, generate=CFA.Model.B)
```

```
plotPowerFitNonNested(Output.B.A, Output.B.B, dat1Mod1=Output.A.A, dat1Mod2=Output.A.B)
plotPowerFitNonNested(Output.B.A, Output.B.B, cutoff=c(AIC=0, BIC=0))

## End(Not run)
```

| | |
|--------------|---|
| plotPowerSig | <i>Plot multiple logistic curves given a significance result matrix</i> |
|--------------|---|

Description

This function will plot the significance results given the value of predictors.

Usage

```
plotPowerSig(sig, x = NULL, xval=NULL, mainName = NULL, useContour = TRUE)
```

Arguments

| | |
|------------|--|
| sig | The data.frame of a significance result, which contains only TRUE for significance and FALSE for not significance. |
| x | The data.frame of the predictor values. The number of rows of the x argument should be equal to the number of rows in the object. |
| xval | The values of predictor that researchers would like to find the fit indices cutoffs from. |
| mainName | A vector of the titles of the graphs |
| useContour | If there are two of sample size, percent completely at random, and percent missing at random are varying, the plotCutoff function will provide 3D graph. Contour graph is a default. However, if this is specified as FALSE, perspective plot is used. |

Value

NONE. Only plot the fit indices distributions.

Author(s)

Sunthud Pornprasertmanit (University of Kansas; <psunthud@ku.edu>)

See Also

- [plotPower](#)

Examples

```
# No example
```

| | |
|-----------|--|
| plotQtile | <i>Build a scatterplot with overlaying line of quantiles of predicted values</i> |
|-----------|--|

Description

Build a scatterplot with overlaying line of quantiles of predicted values

Usage

```
plotQtile(x, y, df=0, qtile=NULL, ...)
```

Arguments

| | |
|-------|--|
| x | The values of the independent variable (e.g., a vector of sample size) |
| y | The values of the dependent variable |
| df | The degree of freedom in spline method |
| qtile | The quantile values used to plot a graph |
| ... | Other arguments in the plot command |

Value

None. This function will plot only.

Author(s)

Sunthud Pornprasertmanit (University of Kansas; psunthud@ku.edu)

Examples

```
# No example
```

| | |
|-------------|---|
| plotScatter | <i>Plot overlaying scatter plots visualizing the power of rejecting misspecified models</i> |
|-------------|---|

Description

This function will find the fit indices cutoff values if not specified and then plot the fit indices value against the value of predictors. The plot will include the fit indices value of the alternative models, the fit indices value of the null model (if specified), and the fit indices cutoffs.

Usage

```
plotScatter(altObject, nullObject=NULL, cutoff=NULL, usedFit = NULL, x, alpha=0.05, df=5)
```

Arguments

| | |
|------------|--|
| altObject | The result object (data.frame) saves the simulation result of fitting the hypothesized model when the hypothesized model is FALSE. |
| nullObject | The result object (data.frame) saves the simulation result of fitting the hypothesized model when the hypothesized model is TRUE. This argument may be not specified if the cutoff is specified. |
| cutoff | A vector of priori cutoffs for fit indices. |
| usedFit | Vector of names of fit indices that researchers wish to plot. |
| x | The data.frame of the predictor values. The number of rows of the x argument should be equal to the number of rows in the object. |
| alpha | A priori alpha level |
| df | The degree of freedom used in spline method in predicting the fit indices by the predictors. If df is 0, the spline method will not be applied. |

Value

NONE. Only plot the fit indices distributions.

Author(s)

Sunthud Pornprasertmanit (University of Kansas; <psunthud@ku.edu>)

See Also

- [plotPowerFit](#)

Examples

```
# No example
```

| | |
|----------------|---|
| popDiscrepancy | <i>Find the discrepancy value between two means and covariance matrices</i> |
|----------------|---|

Description

Find the discrepancy value between two means and covariance matrices

Usage

```
popDiscrepancy(paramM, paramCM, misspecM, misspecCM)
```

Arguments

| | |
|-----------|---|
| paramM | The model-implied mean from the real parameters |
| paramCM | The model-implied covariance matrix from the real parameters |
| misspecM | The model-implied mean from the real and misspecified parameters |
| misspecCM | The model-implied covariance matrix from the real and misspecified parameters |

Details

The discrepancy value (F_0 ; Browne & Cudeck, 1992) is calculated by

$$F_0 = \text{tr}(\tilde{\Sigma}\Sigma^{-1}) - \log|\tilde{\Sigma}\Sigma^{-1}| - p + (\tilde{\mu} - \mu)' \Sigma^{-1} (\tilde{\mu} - \mu).$$

where μ is the model-implied mean from the real parameters, Σ is the model-implied covariance matrix from the real parameters, $\tilde{\mu}$ is the model-implied mean from the real and misspecified parameters, $\tilde{\Sigma}$ is the model-implied covariance matrix from the real and misspecified parameter, p is the number of indicators.

Value

The discrepancy between two means and covariance matrices

Author(s)

Sunthud Pornprasertmanit (University of Kansas; <psunthud@ku.edu>)

References

Browne, M. W., & Cudeck, R. (1992). Alternative ways of assessing model fit. *Sociological Methods & Research*, 21, 230-258.

Examples

```
m1 <- rep(0, 3)
m2 <- c(0.1, -0.1, 0.05)
S1 <- matrix(c(1, 0.6, 0.5, 0.6, 1, 0.4, 0.5, 0.4, 1), 3, 3)
S2 <- matrix(c(1, 0.55, 0.55, 0.55, 1, 0.55, 0.55, 0.55, 1), 3, 3)
popDiscrepancy(m1, S1, m2, S2)
```

popMisfitMACS

Find population misfit by sufficient statistics

Description

Find the value quantifying the amount of population misfit: F_0 , RMSEA, and SRMR.

Usage

```
popMisfitMACS(paramM, paramCM, misspecM, misspecCM, dfParam=NULL, fit.measures="all")
```

Arguments

| | |
|--------------|--|
| paramM | The model-implied mean from the real parameters |
| paramCM | The model-implied covariance matrix from the real parameters |
| misspecM | The model-implied mean from the real and misspecified parameters |
| misspecCM | The model-implied covariance matrix from the real and misspecified parameters |
| dfParam | The degree of freedom of the real model |
| fit.measures | The names of indices used to calculate population misfit. There are three types of misfit: 1) discrepancy function ("f0"; see popDiscrepancy), 2) root mean squared error of approximation ("rmsea"; Equation 12 in Browne & Cudeck, 1992), and 3) standardized root mean squared residual ("srmr") |

Details

The root mean squared error of approximation (RMSEA) is calculated by

$$RMSEA = \sqrt{\frac{F_0}{df}}$$

where F_0 is the discrepancy value between two means vectors and covariance matrices (see [popDiscrepancy](#)) and df is the degree of freedom in the real model.

The standardized root mean squared residual can be calculated by

$$SRMR = \sqrt{\frac{2 \sum_i \sum_{j \leq i} \left(\frac{s_{ij}}{\sqrt{s_{ii}} \sqrt{s_{jj}}} - \frac{\hat{\sigma}_{ij}}{\sqrt{\hat{\sigma}_{ii}} \sqrt{\hat{\sigma}_{jj}}} \right)}{p(p+1)}}$$

where s_{ij} is the observed covariance between indicators i and j , $\hat{\sigma}_{ij}$ is the model-implied covariance between indicators i and j , p is the number of indicators.

Value

The vector of the misfit indices

Author(s)

Sunthud Pornprasertmanit (University of Kansas; <psunthud@ku.edu>)

References

Browne, M. W., & Cudeck, R. (1992). Alternative ways of assessing model fit. *Sociological Methods & Research*, 21, 230-258.

Examples

```
m1 <- rep(0, 3)
m2 <- c(0.1, -0.1, 0.05)
S1 <- matrix(c(1, 0.6, 0.5, 0.6, 1, 0.4, 0.5, 0.4, 1), 3, 3)
S2 <- matrix(c(1, 0.55, 0.55, 0.55, 1, 0.55, 0.55, 0.55, 1), 3, 3)
popMisfitMACS(m1, S1, m2, S2)
```

predProb

Function to get predicted probabilities from logistic regression

Description

Function to get predicted probabilities from logistic regression

Usage

```
predProb(newdat, glmObj)
```

Arguments

| | |
|--------|--|
| newdat | A vector of values for all predictors, including the intercept |
| glmObj | An object from a fitted glm run with a logit link |

Value

Predictive probability of success given the values in the newdat argument.

Author(s)

Alexander M. Schoemann (University of Kansas; <schoemann@ku.edu>)

See Also

- [continuousPower](#)
- [getPower](#)

Examples

```
# No example
```

| | |
|----------------|---|
| printIfNotNull | <i>Provide basic summary of each object if that object is not NULL.</i> |
|----------------|---|

Description

Provide basic summary of each object if that object is not NULL. This function is mainly used in the summary function from the linkS4class{SimSet} object.

Usage

```
printIfNotNull(object, name=NULL)
```

Arguments

| | |
|--------|--|
| object | The target object to be printed, which can be linkS4class{SimMatrix}, linkS4class{SymMatrix}, or linkS4class{SimVector}. |
| name | The name of the target object |

Value

None. This function will print only.

Author(s)

Sunthud Pornprasertmanit (University of Kansas; psunthud@ku.edu)

Examples

```
# This function is not public

# AL <- simVector(rep(NA, 5), "0")
# printIfNotNull(AL, "Factor mean")
```

| | |
|--------|---------------------------------------|
| pValue | <i>Find p-values (1 - percentile)</i> |
|--------|---------------------------------------|

Description

This function will provide p value from comparing number and vector or the analytic result to the observed data (in [lavaan](#)) and the simulation result (in [SimResult](#)).

Usage

```
pValue(target, dist, ...)
```

Arguments

| | |
|--------|---|
| target | A value, multiple values, or a model output object used to find p values. This argument could be a cutoff of a fit index. |
| dist | The comparison distribution, which can be a vector of numbers, a data frame, or a result object. |
| ... | Other values that will be explained specifically for each class |

Details

In comparing fit indices, the p value is the proportion of the number of replications that provide poorer fit (e.g., less CFI value or greater RMSEA value) than the analysis result from the observed data. If the target is a critical value (e.g., fit index cutoff) and the dist is the sampling distribution underlying the alternative hypothesis, this function can provide a statistical power.

Value

Mostly, this function provides a vector of p values based on the comparison. If the target is a model output object and dist is a result object, the p values of fit indices are provided, as well as two additional values: andRule and orRule. The andRule is based on the principle that the model is retained only when all fit indices provide good fit. The proportion is calculated from the number of replications that have all fit indices indicating a better model than the observed data. The proportion from the andRule is the most stringent rule in retaining a hypothesized model. The orRule is based on the principle that the model is retained only when at least one fit index provides good fit. The proportion is calculated from the number of replications that have at least one fit index indicating a better model than the observed data. The proportion from the orRule is the most lenient rule in retaining a hypothesized model.

Methods

signature(target="numeric", dist="vector") This method will find the p value (quantile rank) of the target value on the dist vector. The additional arguments are revDirec, x, xval, condCutoff, and df. The revDirec is a logical argument whether to reverse the direction of comparison. If TRUE, the proportion of the dist that is lower than target value is reported. If FALSE, the proportion of the dist that is higher than the target value is reported. The x is the data.frame of the predictor values. The number of rows of the x argument should be equal to the number of rows in the dist. The xval is the values of predictor that researchers would like to find the fit indices cutoffs from. The condCutoff is a logical argument. If TRUE, the cutoff is applicable only a given value of xval. If FALSE, the cutoff is applicable in any

values of predictor. The *df* is the degree of freedom used in spline method in predicting the fit indices by the predictors. If *df* is 0, the spline method will not be applied.

signature(target="numeric", dist="data.frame") This method will find the *p* value of each columns in the *dist* based on the value specified in the *target*. The additional arguments are *revDirec*, *x*, *xval*, *df*, and *asLogical*. The *revDirec* is a logical vector whether to reverse the direction of comparison. If TRUE, the proportion of the *dist* that is lower than *target* value is reported. If FALSE, the proportion of the *dist* that is higher than the *target* value is reported. The *x* is the *data.frame* of the predictor values. The number of rows of the *x* argument should be equal to the number of rows in the *dist*. The *xval* is the values of predictor that researchers would like to find the fit indices cutoffs from. The *df* is the degree of freedom used in spline method in predicting the fit indices by the predictors. If *df* is 0, the spline method will not be applied. The *asLogical* is to provide the result as the matrix of significance result (TRUE) or just the proportion of significance result (FALSE).

Author(s)

Sunthud Pornprasertmanit (University of Kansas; <psunthud@ku.edu>)

See Also

- [SimResult](#) to run a simulation study

Examples

```
## Not run:
# Compare number with a vector
pValue(0.5, rnorm(1000, 0, 1))

# Compare numbers with a data frame
pValue(c(0.5, 0.2), data.frame(rnorm(1000, 0, 1), runif(1000, 0, 1)))

# Compare an analysis result with a result of simulation study
library(lavaan)
loading <- matrix(0, 9, 3)
loading[1:3, 1] <- NA
loading[4:6, 2] <- NA
loading[7:9, 3] <- NA
targetmodel <- estmodel(LY=loading, modelType="CFA", indLab=paste("x", 1:9, sep=""))
out <- analyze(targetmodel, HolzingerSwineford1939)

loading.trivial <- matrix("runif(1, -0.2, 0.2)", 9, 3)
loading.trivial[is.na(loading)] <- 0
mismodel <- model.lavaan(out, std=TRUE, LY=loading.trivial)

simout <- sim(20, n=nrow(HolzingerSwineford1939), mismodel)
pValue(out, simout)

## End(Not run)
```

| | |
|------------------|---|
| pValueCondCutoff | <i>Find a p value when the target is conditional (valid) on a specific value of a predictor</i> |
|------------------|---|

Description

Find a p value when the target is conditional (valid) on a specific value of a predictor. That is, the target value is applicable only a given value of a predictor.

Usage

```
pValueCondCutoff(target, dist, revDirec = FALSE, x = NULL, xval = NULL, df = 0)
```

Arguments

| | |
|----------|--|
| target | A target value used to find p values. |
| dist | The comparison distribution, which can be a vector of numbers, a data frame, or a result object. |
| revDirec | A logical argument whether to reverse the direction of comparison. If TRUE, the proportion of the dist that is lower than target value is reported. If FALSE, the proportion of the dist that is higher than the target value is reported. |
| x | the data.frame of the predictor values. The number of rows of the x argument should be equal to the number of rows in the dist |
| xval | the values of predictor that researchers would like to find the fit indices cutoffs from. |
| df | the degree of freedom used in spline method in predicting the fit indices by the predictors. If df is 0, the spline method will not be applied. |

Value

A vector of p values based on the comparison.

Author(s)

Sunthud Pornprasertmanit (University of Kansas; <psunthud@ku.edu>)

See Also

- [pValue](#)

Examples

```
# No example
```

pValueNested

*Find p-values (1 - percentile) for a nested model comparison***Description**

This function will provide p value from comparing the differences in fit indices between nested models with the simulation results of both parent and nested models when the nested model is true.

Usage

```
pValueNested(outNested, outParent, simNested, simParent, usedFit = NULL,
nVal = NULL, pmMCARval = NULL, pmMARval = NULL, df = 0)
```

Arguments

| | |
|-----------|---|
| outNested | lavaan that saves the analysis result of the nested model from the target dataset |
| outParent | lavaan that saves the analysis result of the parent model from the target dataset |
| simNested | SimResult that saves the analysis results of nested model from multiple replications |
| simParent | SimResult that saves the analysis results of parent model from multiple replications |
| usedFit | Vector of names of fit indices that researchers wish to getCutoffs from. The default is to getCutoffs of all fit indices. |
| nVal | The sample size value that researchers wish to find the p value from. |
| pmMCARval | The percent missing completely at random value that researchers wish to find the p value from. |
| pmMARval | The percent missing at random value that researchers wish to find the the p value from. |
| df | The degree of freedom used in spline method in predicting the fit indices by the predictors. If df is 0, the spline method will not be applied. |

Details

In comparing fit indices, the p value is the proportion of the number of replications that provide less preference for nested model (e.g., larger negative difference in CFI values or larger positive difference in RMSEA values) than the analysis result from the observed data.

Value

This function provides a vector of p values based on the comparison of the difference in fit indices from the real data with the simulation result. The p values of fit indices are provided, as well as two additional values: andRule and orRule. The andRule is based on the principle that the model is retained only when all fit indices provide good fit. The proportion is calculated from the number of replications that have all fit indices indicating a better model than the observed data. The proportion from the andRule is the most stringent rule in retaining a hypothesized model. The orRule is based on the principle that the model is retained only when at least one fit index provides good fit. The proportion is calculated from the number of replications that have at least one fit index indicating a better model than the observed data. The proportion from the orRule is the most lenient rule in retaining a hypothesized model.

Author(s)

Sunthud Pornprasertmanit (University of Kansas; <psunthud@ku.edu>)

See Also

- [SimResult](#) to run a simulation study

Examples

```
## Not run:
library(lavaan)

LY <- matrix(1, 4, 2)
LY[,2] <- 0:3
PS <- matrix(NA, 2, 2)
TY <- rep(0, 4)
AL <- rep(NA, 2)
TE <- diag(NA, 4)
nested <- estmodel(LY=LY, PS=PS, TY=TY, AL=AL, TE=TE, modelType="CFA", indLab=paste("t", 1:4, sep=""))

LY2 <- matrix(1, 4, 2)
LY2[,2] <- c(0, NA, NA, 3)
parent <- estmodel(LY=LY2, PS=PS, TY=TY, AL=AL, TE=TE, modelType="CFA", indLab=paste("t", 1:4, sep=""))

outNested <- analyze(nested, Demo.growth)
outParent <- analyze(parent, Demo.growth)

loadingMis <- matrix(0, 4, 2)
loadingMis[2:3, 2] <- "runif(1, -0.1, 0.1)"
datamodel <- model.lavaan(outNested, LY=loadingMis)

n <- nrow(Demo.growth)

simNestedNested <- sim(30, n=n, nested, generate=datamodel)
simNestedParent <- sim(30, n=n, parent, generate=datamodel)

pValueNested(outNested, outParent, simNestedNested, simNestedParent)

## End(Not run)
```

pValueNonNested

Find p-values (1 - percentile) for a non-nested model comparison

Description

This function will provide p value from comparing the results of fitting real data into two models against the simulation from fitting the simulated data from both models into both models. The p values from both sampling distribution under the datasets from the first and the second models are reported.

Usage

```
pValueNonNested(outMod1, outMod2, dat1Mod1, dat1Mod2, dat2Mod1, dat2Mod2,
  usedFit = NULL, nVal = NULL, pmMCArval = NULL, pmMARval = NULL, df = 0,
  onetailed=FALSE)
```

Arguments

| | |
|-----------|---|
| outMod1 | lavaan that saves the analysis result of the first model from the target dataset |
| outMod2 | lavaan that saves the analysis result of the second model from the target dataset |
| dat1Mod1 | SimResult that saves the simulation of analyzing Model 1 by datasets created from Model 1 |
| dat1Mod2 | SimResult that saves the simulation of analyzing Model 2 by datasets created from Model 1 |
| dat2Mod1 | SimResult that saves the simulation of analyzing Model 1 by datasets created from Model 2 |
| dat2Mod2 | SimResult that saves the simulation of analyzing Model 2 by datasets created from Model 2 |
| usedFit | Vector of names of fit indices that researchers wish to getCutoffs from. The default is to getCutoffs of all fit indices. |
| nVal | The sample size value that researchers wish to find the p value from. |
| pmCARval | The percent missing completely at random value that researchers wish to find the p value from. |
| pmMARval | The percent missing at random value that researchers wish to find the the p value from. |
| df | The degree of freedom used in spline method in predicting the fit indices by the predictors. If df is 0, the spline method will not be applied. |
| onetailed | If TRUE, the function will convert the p value based on two-tailed test. |

Details

In comparing fit indices, the p value is the proportion of the number of replications that provide less preference for either model 1 or model 2 than the analysis result from the observed data. In two-tailed test, the function will report the proportion of values under the sampling distribution that are more extreme than one obtained from real data. If the resulting p value is high ($> .05$) on one model and low ($< .05$) in the other model, the model with high p value is preferred. If the p values are both high or both low, the decision is undetermined.

Value

This function provides a vector of p values based on the comparison of the difference in fit indices from the real data with the simulation results. The p values of fit indices are provided, as well as two additional values: andRule and orRule. The andRule is based on the principle that the model is retained only when all fit indices provide good fit. The proportion is calculated from the number of replications that have all fit indices indicating a better model than the observed data. The proportion from the andRule is the most stringent rule in retaining a hypothesized model. The orRule is based on the principle that the model is retained only when at least one fit index provides good fit. The proportion is calculated from the number of replications that have at least one fit index indicating a better model than the observed data. The proportion from the orRule is the most lenient rule in retaining a hypothesized model.

Author(s)

Sunthud Pornprasertmanit (University of Kansas; <psunthud@ku.edu>)

See Also

- [SimResult](#) to run a simulation study

Examples

```
## Not run:
library(lavaan)
loading <- matrix(0, 11, 3)
loading[1:3, 1] <- NA
loading[4:7, 2] <- NA
loading[8:11, 3] <- NA
path.A <- matrix(0, 3, 3)
path.A[2:3, 1] <- NA
path.A[3, 2] <- NA
model.A <- estmodel(LY=loading, BE=path.A, modelType="SEM", indLab=c(paste("x", 1:3, sep=""), paste("y", 1:3, sep="")))

out.A <- analyze(model.A, PoliticalDemocracy)

path.B <- matrix(0, 3, 3)
path.B[1:2, 3] <- NA
path.B[1, 2] <- NA
model.B <- estmodel(LY=loading, BE=path.B, modelType="SEM", indLab=c(paste("x", 1:3, sep=""), paste("y", 1:3, sep="")))

out.B <- analyze(model.B, PoliticalDemocracy)

loading.mis <- matrix("runif(1, -0.2, 0.2)", 11, 3)
loading.mis[is.na(loading)] <- 0

datamodel.A <- model.lavaan(out.A, std=TRUE, LY=loading.mis)
datamodel.B <- model.lavaan(out.B, std=TRUE, LY=loading.mis)

n <- nrow(PoliticalDemocracy)

output.A.A <- sim(5, n=n, model.A, generate=datamodel.A)
output.A.B <- sim(5, n=n, model.B, generate=datamodel.A)
output.B.A <- sim(5, n=n, model.A, generate=datamodel.B)
output.B.B <- sim(5, n=n, model.B, generate=datamodel.B)

# The output may contain some warnings here. When the number of replications increases (e.g., 1000), the warning
# pValueNonNested(out.A, out.B, output.A.A, output.A.B, output.B.A, output.B.B) will appear.

## End(Not run)
```

pValueVariedCutoff

Find a p value when the cutoff is specified as a vector given the values of predictors

Description

Find a p value when the cutoff is specified as a vector given the values of predictors.

Usage

```
pValueVariedCutoff(cutoff, obtainedValue, revDirec = FALSE, x = NULL, xval = NULL)
```

Arguments

| | |
|---------------|--|
| cutoff | A vector of values used to find p values. Each value in the vector should be the target value conditional (applicable) to each value of the predictors (x) respectively. |
| obtainedValue | The comparison distribution, which can be a vector of numbers, a data frame, or a result object. |
| revDirec | A logical argument whether to reverse the direction of comparison. If TRUE, the proportion of the dist that is lower than target value is reported. If FALSE, the proportion of the dist that is higher than the target value is reported. |
| x | the data.frame of the predictor values. The number of rows of the x argument should be equal to the number of rows in the dist |
| xval | the values of predictor that researchers would like to find the fit indices cutoffs from. |

Value

A vector of p values based on the comparison.

Author(s)

Sunthud Pornprasertmanit (University of Kansas; <psunthud@ku.edu>)

See Also

- [pValue](#)

Examples

```
# No example
```

revText

Reverse the proportion value by subtracting it from 1

Description

Reverse the proportion value by subtracting it from 1. This function can reverse a value reported in text, such as from "> .98" to "< .02"

Usage

```
revText(val)
```

Arguments

| | |
|-----|--------------------------|
| val | The value to be reversed |
|-----|--------------------------|

Value

The reversed value or text

Author(s)

Sunthud Pornprasertmanit (University of Kansas; <psunthud@ku.edu>)

Examples

```
# This is a private function.

# revText(.96)
# revText("> .60")
```

setPopulation

Set the data generation population model underlying an object

Description

This function will set the data generation population model to be an appropriate one. If the appropriate data generation model is put (the same model as the analysis model), the additional features can be seen when we run a [summary](#) function on the target object, such as bias in parameter estimates or percentage coverage.

Usage

```
setPopulation(target, population)
```

Arguments

| | |
|------------|---|
| target | The result object that you wish to set the data generation population model (linkS4class{SimResult}). |
| population | The population parameters specified in the linkS4class{SimSem} object |

Value

The target object that is changed the parameter.

Author(s)

Sunthud Pornprasertmanit (University of Kansas; <psunthud@ku.edu>)

See Also

- [SimResult](#) for result object

Examples

```
# See each class for an example.
## Not run:

loading <- matrix(0, 7, 3)
loading[1:3, 1] <- NA
loading[4:6, 2] <- NA
loading[1:7, 3] <- NA
loadingVal <- matrix(0, 7, 3)
```



```

loadingVal[1:3, 1] <- "runif(1, 0.5, 0.7)"
loadingVal[4:6, 2] <- "runif(1, 0.5, 0.7)"
loadingVal[1:6, 3] <- "runif(1, 0.3, 0.5)"
loadingVal[7, 3] <- 1
loading.mis <- matrix("runif(1, -0.2, 0.2)", 7, 3)
loading.mis[is.na(loading)] <- 0
loading.mis[,3] <- 0
loading.mis[7,] <- 0
LY <- bind(loading, loadingVal, misspec=loading.mis)

RPS <- binds(diag(3))

path <- matrix(0, 3, 3)
path[2, 1] <- NA
BE <- bind(path, "runif(1, 0.3, 0.5)")

RTE <- binds(diag(7))

VY <- bind(c(rep(NA, 6), 0), c(rep(1, 6), ""))

datamodel <- model(LY=LY, RPS=RPS, BE=BE, RTE=RTE, VY=VY, modelType="SEM")

loading <- matrix(0, 7, 3)
loading[1:3, 1] <- NA
loading[4:6, 2] <- NA
loading[7, 3] <- NA
path <- matrix(0, 3, 3)
path[2, 1] <- NA
path[1, 3] <- NA
path[2, 3] <- NA
errorCov <- diag(NA, 7)
errorCov[7, 7] <- 0
facCov <- diag(3)
analysis <- estmodel(LY=loading, BE=path, TE=errorCov, PS=facCov, modelType="SEM", indLab=paste("y", 1:7, s

Output <- sim(100, n=200, analysis, generate=datamodel)

loadingVal <- matrix(0, 7, 3)
loadingVal[1:3, 1] <- 0.6
loadingVal[4:6, 2] <- 0.6
loadingVal[7, 3] <- 1
LY <- bind(loading, loadingVal)
pathVal <- matrix(0, 3, 3)
pathVal[2, 1] <- 0.4
pathVal[1, 3] <- 0.4
pathVal[2, 3] <- 0.4
BE <- bind(path, pathVal)
PS <- binds(facCov)
errorCovVal <- diag(0.64, 7)
errorCovVal[7, 7] <- 0
TE <- binds(errorCov, errorCovVal)
population <- model(LY=LY, PS=PS, BE=BE, TE=TE, modelType="SEM")
Output <- setPopulation(Output, population)
summary(Output)

## End(Not run)

```

sim

*Run a monte carlo simulation with a structural equation model.***Description**

TBA

Usage

```
sim(nRep, model, n, generate = NULL, rawData = NULL, miss = NULL, datafun=NULL, outfun=NULL,
pmMCAR = NULL, pmMAR = NULL, facDist = NULL, indDist = NULL, errorDist = NULL, sequential = FALSE,
modelBoot = FALSE, realData = NULL, maxDraw = 50, misfitType = "f0",
misfitBounds = NULL, averageNumMisspec = NULL, optMisfit=NULL, optDraws = 50,
aux = NULL, seed = 123321, silent = FALSE, multicore = FALSE, cluster = FALSE, numProc = NULL,
paramOnly = FALSE, dataOnly=FALSE, ...)
```

Arguments

| | |
|----------|---|
| nRep | Number of replications. Users can specify as NULL and specify n, pmMCAR, and pmMAR |
| model | SimSem object created by model . Will be used to generate data and analyze it. |
| n | Sample size. This argument is not necessary except the user wish to vary sample size across replications. The sample size here is a vector of sample size in integers. For the random distribution object, if the resulting value has decimal, the value will be rounded. |
| generate | SimSem object created by model . If included, this will be used to generate data instead of |
| rawData | If included, this data is used for simulations instead of being generated from the SimSem template. Should be a list of length nRep. |
| miss | Missing data handling template, created by the function miss . |
| datafun | Function to be applied to generated data set at each replication. |
| outfun | Function to be applied to the lavaan output at each replication. The output of this function in each replication will be saved in the simulation output (SimResult). The extra outputs can be obtained by the getExtraOutput function. |
| pmMCAR | The percent completely missing at random. This argument is not necessary except the user wish to vary percent missing completely at random across replications. The pmMCAR here is a vector of percent missing, which the values can be in between 0 and 1 only. The specification of objMissing is not needed (but is needed if users wish to specify complex missing value data generation or wish to use multiple imputation). |
| pmMAR | The percent missing at random. This argument is not necessary except the user wish to vary percent missing at random across replications. The pmMAR here is a vector of percent missing, which the values can be in between 0 and 1 only. The specification of objMissing is not needed (but is needed if users wish to specify complex missing value data generation or wish to use multiple imputation). |
| facDist | A SimDataDist object or list of objects for the distribution of factors. If one object is passed, all factors will have the same distribution. Use when sequential is TRUE. |

| | |
|-------------------|---|
| indDist | A SimDataDist object or list of objects for a distribution of indicators. If one object is passed, each indicator will have the same distribution. Use when sequential is FALSE. |
| errorDist | An object or list of objects of type SimDataDist indicating the distribution of errors. If a single SimDataDist is specified, each error will be generated with that distribution. |
| sequential | If TRUE, use a sequential method to create data such that the data from factor are generated first and apply to a set of equations to obtain the data of indicators. If FALSE, create data directly from model-implied mean and covariance of indicators. |
| modelBoot | When specified, a model-based bootstrap is used for data generation. See draw for further information. This argument requires real data to be passed to <code>realData</code> . |
| realData | A <code>data.frame</code> containing real data. The data generated will follow the distribution of this data set. |
| maxDraw | Integer specifying the maximum number of attempts to draw a valid set of parameters (no negative error variance, standardized coefficients over 1). |
| misfitType | Character vector indicating the fit measure used to assess the misfit of a set of parameters. Can be "f0", "rmsea", "srmr", or "all". |
| misfitBounds | Vector that contains upper and lower bounds of the misfit measure. Sets of parameters drawn that are not within these bounds are rejected. |
| averageNumMisspec | TRUE or FALSE. ?? |
| optMisfit | Character vector of either "min" or "max" indicating either maximum or minimum optimized misfit. If not null, the set of parameters out of the number of draws in "optDraws" that has either the maximum or minimum misfit of the given misfit type will be returned. |
| optDraws | Number of parameter sets to draw if <code>optMisfit</code> is not null. The set of parameters with the maximum or minimum misfit will be returned. |
| aux | The names of auxiliary variables saved in a vector |
| seed | Random number seed. Reproducibility across multiple cores or clusters is ensured using R's <code>Lecuyer</code> package. |
| silent | If TRUE, suppress warnings. |
| multicore | Use multiple processors within a computer. Specify as TRUE to use it. |
| cluster | Not applicable now. Use for specify nodes in <code>hpc</code> in order to be parallelizable. |
| numProc | Number of processors for using multiple processors. If it is NULL, the package will find the maximum number of processors. |
| paramOnly | If TRUE, only the parameters from each replication will be returned. |
| dataOnly | If TRUE, only the data generated from each replication will be returned. |
| ... | Additional arguments to be passed to <code>lavaan</code> . |

Value

TBA

Author(s)

Patrick Miller (Univeristy of Notre Dame; <pmille13@nd.edu>) Sunthud Pornprasertmanit (University of Kansas; <psunthud@ku.edu>)

Examples

```

loading <- matrix(0, 6, 2)
loading[1:3, 1] <- NA
loading[4:6, 2] <- NA
LY <- bind(loading, 0.7)

latent.cor <- matrix(NA, 2, 2)
diag(latent.cor) <- 1
RPS <- binds(latent.cor, 0.5)

RTE <- binds(diag(6))

VY <- bind(rep(NA,6),2)

CFA.Model <- model(LY = LY, RPS = RPS, RTE = RTE, modelType = "CFA")

Output <- sim(20, CFA.Model,n=200)
summary(Output)

```

| | |
|-------------------|---------------------|
| SimDataDist-class | Class "SimDataDist" |
|-------------------|---------------------|

Description

This class will provide the distribution of a dataset.

Objects from the Class

Objects can be created by `bindDist` function. It can also be called from the form `new("SimDataDist", ...)`.

Slots

p: Number of variables

margins: A character vector specifying all the marginal distributions

paramMargins: A list whose each component is a list of named components, giving the parameter values of the marginal distributions.

keepScale: Transform back to retain the mean and standard deviation of a variable equal to the model implied mean and standard deviation (with sampling error)

reverse: To mirror each variable or not. If TRUE, reverse the distribution of a variable (e.g., from positive skewed to negative skewed).

Methods

- `summaryTo` summarize the object
- `plotDistTo` plot a density distribution (for one variable) or a contour plot (for two variables). If the object has more than two variables, the `var` argument can be used to select the index of plotting variables. For two variables, the default is to have correlation of 0. To change a correlation, the `r` argument can be used. The `xlim` and `ylim` can be specified to set the ranges of variables.
- `extractExtract` elements from an object. The next argument is the position of the object to be extracted.

Author(s)

Sunthud Pornprasertmanit (University of Kansas; <psunthud@ku.edu>)

See Also

- [bindDist](#) The constructor of this class.

Examples

```
# Need to be fixed

showClass("SimDataDist")

d1 <- list(df=2)
d2 <- list(df=3)
d3 <- list(df=4)
d4 <- list(df=5)
d5 <- list(df=3)
d6 <- list(df=4)
d7 <- list(df=5)
d8 <- list(df=6)

dist <- bindDist(c(rep("t", 4), rep("chisq", 8)), d1, d2, d3, d4, d5, d6, d7, d8, d5, d6, d7, d8)
```

SimMatrix-class

Matrix object: Random parameters matrix

Description

This object can be used to represent a matrix in SEM model. It contains free parameters, fixed values, and starting values. This object can be represented factor loading matrix or regression coefficient matrix.

Objects from the Class

This object is created by [bind](#) function. Objects can be also created by calls of the form `new("SimMatrix", ...)`.

Slots

free: TBA
 popParam: TBA
 misspec: TBA
 symmetric: TBA

Methods

[summaryShort](#) Provides a short summary of all information in the object
[summary](#) Provides a thorough description of all information in the object

Author(s)

Sunthud Pornprasertmanit (University of Kansas; <psunthud@ku.edu>)

See Also

- [SimVector](#) for random parameter vector.

Examples

```
showClass("SimMatrix")

loading <- matrix(0, 6, 2)
loading[1:3, 1] <- NA
loading[4:6, 2] <- NA
loadingValues <- matrix(0, 6, 2)
loadingValues[1:3, 1] <- 0.7
loadingValues[4:6, 2] <- 0.7
LX <- bind(loading, loadingValues)
summary(LX)
# run(LX)

LY <- bind(loading, "rnorm(1, 0.6, 0.05)")
summary(LY)
# run(LY)
```

| | |
|------------------|---------------------------|
| SimMissing-class | <i>Class "SimMissing"</i> |
|------------------|---------------------------|

Description

Missing information imposing on the complete dataset

Objects from the Class

Objects can be created by `simMissing` function. It can also be called from the form `new("SimMissing", ...)`.

Slots

cov: Column indices of any normally distributed covariates used in the data set.

pmMCAR: Decimal percent of missingness to introduce completely at random on all variables.

pmMAR: Decimal percent of missingness to introduce using the listed covariates as predictors.

nforms: The number of forms for planned missing data designs, not including the shared form.

itemGroups: List of lists of item groupings for planned missing data forms. Without this, items will be divided into groups sequentially (e.g. 1-3,4-6,7-9,10-12)

twoMethod: Vector of (percent missing, column index). Will put a given percent missing on that column in the matrix to simulate a two method planned missing data research design.

prAttr: Probability (or vector of probabilities) of an entire case being removed due to attrition at a given time point. See [imposeMissing](#) for further details.

package: The package used in multiple imputation. If "default", the full-information maximum likelihood is used.

timePoints: Number of timepoints items were measured over. For longitudinal data, planned missing designs will be implemented within each timepoint.

ignoreCols: The columns not imposed any missing values for any missing data patterns

threshold: The threshold of covariates that divide between the area to impose missing and the area not to impose missing. The default threshold is the mean of the covariate.

covAsAux: If TRUE, the covariate listed in the object will be used as auxiliary variables when putting in the model object. If FALSE, the covariate will be included in the analysis.

logical: A matrix of logical values (TRUE/FALSE). If a value in the dataset is corresponding to the TRUE in the logical matrix, the value will be missing.

args: A list of additional options to be passed to the multiple imputation function in each package.

Methods

- [summary](#) To summarize the object

Author(s)

Patrick Miller(University of Kansas; <patr1ckm@ku.edu>) Alexander M. Schoemann (University of Kansas; <schoemann@ku.edu>) Kyle Lang (University of Kansas; <kylelang@ku.edu>) Sunthud Pornprasertmanit (University of Kansas; <psunthud@ku.edu>)

See Also

- [imposeMissing](#) for directly imposing missingness into a dataset.

Examples

```
# No Example
```

| | |
|-----------------|-------------------|
| SimResult-class | Class "SimResult" |
|-----------------|-------------------|

Description

This class will save data analysis results from multiple replications and ready to find some useful statistics, such as fit indices cutoffs or power.

Objects from the Class

Objects can be created by [sim](#). It can also be called from the form `new("SimResult", ...)`.

Slots

modelType: Analysis model type (CFA, Path, or SEM)

nRep: Number of replications have been created and run simulated data.

coef: Parameter estimates from each replication

se: Standard errors of parameter estimates from each replication

fit: Fit Indices values from each replication

converged: Number of convergence replications

seed: Seed number.

paramValue: Population model underlying each simulated dataset.

misspecValue: Misspecified-parameter values that are imposed on the population model in each replication.

popFit: The amount of population misfit. See details at [popMisfitMACS](#)

FMI1: Fraction Missing Method 1.

FMI2: Fraction Missing Method 2.

stdCoef: Standardized coefficients from each replication

n: Sample size of the analyzed data.

pmMCAR: Percent missing completely at random.

pmMAR: Percent missing at random.

extraOut: Extra outputs obtained from running the function specified in outfun argument.

timing: Time elapsed in each phase of the simulation.

Methods

- [getCutoff](#) to getCutoff of fit indices based on a priori alpha level.
- [getPowerFit](#) to getPowerFit of rejection when the simResult is the alternative hypothesis and users specify cutoffs of the fit indices.
- [plotCutoff](#) to plot null hypothesis sampling distributions of fit indices with an option to draw fit indices cutoffs by specifying a priori alpha level.
- [plotPowerFit](#) to plot alternative hypothesis (and null hypothesis) with a priori cutoffs or alpha level.
- [summary](#) to summarize the result output
- [summaryParam](#) to summarize all parameter estimates
- [anova](#) find the averages of model fit statistics and indices for nested models, as well as the differences of model fit indices among models. This function requires at least two SimResult objects. See [anova](#) for further details.
- [summaryPopulation](#) to summarize the data generation population underlying the simulation study.
- [getPopulation](#) to extract the data generation population underlying the simulation study. This method will return a data frame of the population underlying each replication.
- [setPopulation](#) to put the appropriate data generation model into the result object. If the appropriate data generation model is put (the same model as the analysis model), the bias in parameter estimates and standard errors will be able to be calculated by the summary function. The first argument is the result object. The second argument can be either data.frame of the population or SimSet of the population. See the 'modeling with covariate' in the manual for an example.

Author(s)

Sunthud Pornprasertmanit (University of Kansas; <psunthud@ku.edu>)

See Also

- [sim](#) for the constructor of this class

Examples

```

showClass("SimResult")
loading <- matrix(0, 6, 1)
loading[1:6, 1] <- NA
LX <- bind(loading, 0.7)
RPH <- binds(diag(1))
RTD <- binds(diag(6))
CFA.Model <- model(LY = LX, RPS = RPH, RTE = RTD, modelType="CFA")

# We make the examples running only 5 replications to save time.
# In reality, more replications are needed.
Output <- sim(5, n=500, CFA.Model)
summary(Output)
getCutoff(Output, 0.05)
summaryParam(Output)
summaryPopulation(Output)

```

SimSem-class

Class "SimSem"

Description

TBA

Objects from the Class

TBA

Slots

pt: TBA

dgen: TBA

modelType: TBA

groupLab: TBA

Methods**summary** Get the summary of model specification**Author(s)**

Patrick Miller (Univeristy of Notre Dame; <pmille13@nd.edu>), Sunthud Pornprasertmanit (University of Kansas; <psunthud@ku.edu>)

See Also

- Create an object this class by CFA, Path Analysis, or SEM model by [model](#).

Examples

```
showClass("SimSem")

loading <- matrix(0, 6, 2)
loading[1:3, 1] <- NA
loading[4:6, 2] <- NA
loadingValues <- matrix(0, 6, 2)
loadingValues[1:3, 1] <- 0.7
loadingValues[4:6, 2] <- 0.7
LX <- bind(loading, loadingValues)
summary(LX)

latent.cor <- matrix(NA, 2, 2)
diag(latent.cor) <- 1
RPH <- binds(latent.cor, 0.5)

# Error Correlation Object
error.cor <- matrix(0, 6, 6)
diag(error.cor) <- 1
RTD <- binds(error.cor)

CFA.Model <- model(LY = LX, RPS = RPH, RTE = RTD, modelType="CFA")
summary(CFA.Model)
#run(CFA.Model)

#CFA.Model2 <- extract(CFA.Model, y=1:3, e=1)
#summary(CFA.Model2)
```

SimVector-class

Vector object: Random parameters vector

Description

This object can be used to represent a vector in SEM model. It contains free parameters, fixed values, and starting values. This object can be represented mean, intercept, or variance vectors.

Objects from the Class

This object is created by [bind](#) function. Objects can be created by calls of the form `new("SimVector", ...)`.

Slots

free: TBA
 popParam: TBA
 misspec: TBA

Methods

[summaryShort](#) Provides a short summary of all information in the object
[summary](#) Provides a thorough description of all information in the object

Author(s)

Sunthud Pornprasertmanit (University of Kansas; <psunthud@ku.edu>)

See Also

[SimMatrix](#) for random parameter matrix

Examples

```
showClass("SimVector")

factor.mean <- rep(NA, 2)
factor.mean.starting <- c(5, 2)
AL <- bind(factor.mean, factor.mean.starting)
#run(AL)
summary(AL)
summaryShort(AL)
```

sortList

Sort two objects in a list

Description

Sort two objects in a list by swapping the values of both objects so that the first object contains the lower value and the second object contains the larger value

Usage

```
sortList(object)
```

Arguments

object The list with two objects (e.g., vector, matrix)

Value

The sorted list

Author(s)

Sunthud Pornprasertmanit (University of Kansas; <psunthud@ku.edu>)

Examples

```
# No example
```

| | |
|-----------------|---|
| summaryConverge | <i>Provide a comparison between the characteristics of convergent replications and nonconvergent replications</i> |
|-----------------|---|

Description

This function provides a comparison between the characteristics of convergent replications and nonconvergent replications. The comparison includes sample size (if varying), percent missing completely at random (if varying), percent missing at random (if varying), parameter values, misspecified-parameter values (if applicable), and population misfit (if applicable).

Usage

```
summaryConverge(object)
```

Arguments

object [SimResult](#) object being described

Value

A list with the following elements:

- Converged The number of convergent and nonconvergent replications
- n Sample size
- pmMCAR Percent missing completely at random
- pmMAR Percent missing at random
- paramValue Parameter values
- misspecValue Misspecified-parameter values
- popFit Population misfit

Each element will provide the head-to-head comparison between convergent and nonconvergent replications properties.

Author(s)

Sunthud Pornprasertmanit (University of Kansas; <psunthud@ku.edu>)

Examples

```
## Not run:
path.BE <- matrix(0, 4, 4)
path.BE[3, 1:2] <- NA
path.BE[4, 3] <- NA
starting.BE <- matrix("", 4, 4)
starting.BE[3, 1:2] <- "runif(1, 0.3, 0.5)"
starting.BE[4, 3] <- "runif(1, 0.5, 0.7)"
mis.path.BE <- matrix(0, 4, 4)
mis.path.BE[4, 1:2] <- "runif(1, -0.1, 0.1)"
BE <- bind(path.BE, starting.BE, misspec=mis.path.BE)
```

```

residual.error <- diag(4)
residual.error[1,2] <- residual.error[2,1] <- NA
RPS <- binds(residual.error, "rnorm(1, 0.3, 0.1)")

loading <- matrix(0, 12, 4)
loading[1:3, 1] <- NA
loading[4:6, 2] <- NA
loading[7:9, 3] <- NA
loading[10:12, 4] <- NA
mis.loading <- matrix("runif(1, -0.3, 0.3)", 12, 4)
mis.loading[is.na(loading)] <- 0
LY <- bind(loading, "runif(1, 0.7, 0.9)", misspec=mis.loading)

mis.error.cor <- matrix("rnorm(1, 0, 0.1)", 12, 12)
diag(mis.error.cor) <- 0
RTE <- binds(diag(12), misspec=mis.error.cor)

SEM.Model <- model(RPS = RPS, BE = BE, LY=LY, RTE=RTE, modelType="SEM")

n1 <- list(mean = 0, sd = 0.1)
chi5 <- list(df = 5)

facDist <- bindDist(c("chisq", "chisq", "norm", "norm"), chi5, chi5, n1, n1)

dat <- generate(SEM.Model, n=500, sequential=TRUE, facDist=facDist)
out <- analyze(SEM.Model, dat, estimator="mlr")

simOut <- sim(50, n=500, SEM.Model, sequential=TRUE, facDist=facDist, estimator="mlr")
summaryConverge(simOut)

## End(Not run)

```

summaryFit

*Provide summary of model fit across replications***Description**

This function will provide fit index cutoffs for values of alpha, and mean fit index values across all replications.

Usage

```
summaryFit(object, alpha = NULL)
```

Arguments

| | |
|--------|--|
| object | SimResult to be summarized |
| alpha | The alpha level used to find the fit indices cutoff. If there is no varying condition, a vector of different alpha levels can be provided. |

Value

A data frame that provides fit statistics cutoffs and means

When `linkS4class{SimResult}` has fixed simulation parameters the first columns are fit index cutoffs for values of alpha and the last column is the mean fit across all replications. Rows are

- Chi Chi-square fit statistic
- AIC Akaike Information Criterion
- BIC Bayesian Information Criterion
- RMSEA Root Mean Square Error of Approximation
- CFI Comparative Fit Index
- TLI Tucker-Lewis Index
- SRMR Standardized Root Mean Residual

When `linkS4class{SimResult}` has random simulation parameters (sample size or percent missing), columns are the fit indices listed above and rows are values of the random parameter.

See details in [popDiscrepancy](#) and [popMisfitMACS](#)

Author(s)

Alexander M. Schoemann (University of Kansas; <schoemann@ku.edu>) Sunthud Pornprasertmanit (University of Kansas; <psunthud@ku.edu>)

See Also

[SimResult](#) for the result object input

Examples

```
loading <- matrix(0, 6, 1)
loading[1:6, 1] <- NA
LX <- bind(loading, 0.7)
RPH <- binds(diag(1))
RTD <- binds(diag(6))
CFA.Model <- model(LY = LX, RPS = RPH, RTE = RTD, modelType="CFA")

# We make the examples running only 5 replications to save time.
# In reality, more replications are needed.
Output <- sim(5, n=500, CFA.Model)
summaryFit(Output)
```

summaryMisspec

Provide summary of the population misfit and misspecified-parameter values across replications

Description

This function provides the summary of the population misfit and misspecified-parameter values across replications. The summary will be provided for the convergent replications only.

Usage

```
summaryMisspec(object)
```

Arguments

object [SimResult](#) object being described

Value

A data frame that provides the summary of population misfit and misspecified-parameter values imposed on the real parameters

Author(s)

Sunthud Pornprasertmanit (University of Kansas; <psunthud@ku.edu>)

See Also

[SimResult](#) for the object input

Examples

```
## Not run:
path <- matrix(0, 4, 4)
path[3, 1:2] <- NA
path[4, 3] <- NA
pathVal <- matrix("", 4, 4)
pathVal[3, 1:2] <- "runif(1, 0.3, 0.5)"
pathVal[4, 3] <- "runif(1, 0.5, 0.7)"
pathMis <- matrix(0, 4, 4)
pathMis[4, 1:2] <- "runif(1, -0.1, 0.1)"
BE <- bind(path, pathVal, pathMis)

residual.error <- diag(4)
residual.error[1,2] <- residual.error[2,1] <- NA
RPS <- binds(residual.error, "rnorm(1, 0.3, 0.1)")

Path.Model <- model(RPS = RPS, BE = BE, modelType="Path")

# The number of replications in actual analysis should be much more than 5
ParamObject <- sim(5, n=200, Path.Model)

summaryMisspec(ParamObject)

## End(Not run)
```

Description

This function will provide averages of parameter estimates, standard deviations of parameter estimates, averages of standard errors, and power of rejection with a priori alpha level for the null hypothesis of parameters equal 0.

Usage

```
summaryParam(object, alpha = 0.05, detail = FALSE)
```

Arguments

| | |
|--------|--|
| object | SimResult object being described |
| alpha | The alpha level used to find the statistical power of each parameter estimate |
| detail | If TRUE, more details about each parameter estimate are provided, such as relative bias, standardized bias, or relative standard error bias. |

Value

A data frame that provides the statistics described above from all parameters. For using with `linkS4class{SimResult}`, each column means

- `Estimate.Average`: Average of parameter estimates across all replications
- `Estimate.SD`: Standard Deviation of parameter estimates across all replications
- `Average.SE`: Average of standard errors across all replications
- `Power (Not equal 0)`: Proportion of significant replications when testing whether the parameters are different from zero
- `Average.Param`: Parameter values or average values of parameters if random parameters are specified
- `SD.Param`: Standard Deviations of parameters. Appeared only when random parameters are specified.
- `Average.Bias`: The difference between parameter estimates and parameter underlying data
- `SD.Bias`: Standard Deviations of bias across all replications. Appeared only when random parameters are specified. This value is the expected value of average standard error when random parameter are specified.
- `Coverage`: The percentage of (1-alpha)% confidence interval covers parameters underlying the data.
- `Rel.Bias`: Relative Bias, which is $(\text{Estimate.Average} - \text{Average.Param}) / \text{Average.Param}$. Hoogland and Boomsma (1998) proposed that the cutoff of .05 may be used for acceptable relative bias. This option will be available when `detail=TRUE`. This value will not be available when parameter values are very close to 0.
- `Std.Bias`: Standardized Bias, which is $(\text{Estimate.Average} - \text{Average.Param}) / \text{Estimate.SD}$ for fixed parameters and $(\text{Estimate.Average} - \text{Average.Param}) / \text{SD.Bias}$ for random parameters. Collins, Schafer, and Kam (2001) recommended that biases will be only noticeable when standardized bias is greater than 0.4 in magnitude. This option will be available when `detail=TRUE`
- `Rel.SE.Bias`: Relative Bias in standard error, which is $(\text{Average.SE} - \text{Estimate.SD}) / \text{Estimate.SD}$ for fixed parameters and $(\text{Average.SE} - \text{SD.Bias}) / \text{SD.Bias}$ for random parameters. Hoogland and Boomsma (1998) proposed that 0.10 is the acceptable level. This option will be available when `detail=TRUE`

Author(s)

Sunthud Pornprasertmanit (University of Kansas; <psunthud@ku.edu>)

References

- Collins, L. M., Schafer, J. L., & Kam, C. M. (2001). A comparison of inclusive and restrictive strategies in modern missing data procedures. *Psychological Methods*, 6, 330-351.
- Hoogland, J. J., & Boomsma, A. (1998). Robustness studies in covariance structure modeling. *Sociological Methods & Research*, 26, 329-367.

See Also

[SimResult](#) for the object input

Examples

```
showClass("SimResult")
loading <- matrix(0, 6, 1)
loading[1:6, 1] <- NA
LX <- bind(loading, 0.7)
RPH <- binds(diag(1))
RTD <- binds(diag(6))
CFA.Model <- model(LY = LX, RPS = RPH, RTE = RTD, modelType="CFA")

# We make the examples running only 5 replications to save time.
# In reality, more replications are needed.
Output <- sim(5, n=500, CFA.Model)
summaryParam(Output)
summaryParam(Output, detail=TRUE)
```

| | |
|-------------------|--|
| summaryPopulation | <i>Summarize the data generation population model underlying a result object</i> |
|-------------------|--|

Description

Summarize the data generation population model underlying a result object

Usage

```
summaryPopulation(object)
```

Arguments

| | |
|--------|---|
| object | The result object that you wish to extract the data generation population model from (<code>linkS4class{SimResult}</code>). |
|--------|---|

Value

None except using for `linkS4class{SimResult}` which the return value is a `data.frame` of the summary of population model across replications.

Author(s)

Sunthud Pornprasertmanit (University of Kansas; <psunthud@ku.edu>)

See Also

- [SimResult](#) for result object

Examples

```
## Not run:
loading <- matrix(0, 6, 1)
loading[1:6, 1] <- NA
LX <- bind(loading, "runif(1, 0.4, 0.9)")
RPH <- binds(diag(1))
RTD <- binds(diag(6))
CFA.Model <- model(LY = LX, RPS = RPH, RTE = RTD, modelType="CFA")

# We will use only 10 replications to save time.
# In reality, more replications are needed.
Output <- sim(10, n=200, model=CFA.Model)
summaryPopulation(Output)

## End(Not run)
```

summaryShort

Provide short summary of an object.

Description

Provide short summary if it is available. Otherwise, it is an alias for summary.

Usage

```
summaryShort(object, ...)
```

Arguments

| | |
|--------|--------------------------------|
| object | Desired object being described |
| ... | any additional arguments |

Value

NONE. This function will print on screen only.

Author(s)

Sunthud Pornprasertmanit (University of Kansas; <psunthud@ku.edu>)

See Also

This is the list of classes that can use summaryShort method.

- [SimMatrix](#)
- [SimVector](#)

Examples

```
loading <- matrix(0, 6, 2)
loading[1:3, 1] <- NA
loading[4:6, 2] <- NA
loadingValues <- matrix(0, 6, 2)
LX <- bind(loading, "runif(1, 0.8, 0.9)")
summaryShort(LX)
```

| | |
|-----------------|--|
| twoTailedPValue | <i>Find two-tailed p value from one-tailed p value</i> |
|-----------------|--|

Description

Find two-tailed p value from one-tailed p value

Usage

```
twoTailedPValue(vec)
```

Arguments

vec A vector of one-tailed p value.

Value

A vector of two-tailed p value.

Author(s)

Sunthud Pornprasertmanit (University of Kansas; <psunthud@ku.edu>)

Examples

```
# No example
```

| | |
|--------------------|--|
| validateCovariance | <i>Validate whether all elements provides a good covariance matrix</i> |
|--------------------|--|

Description

Validate whether all elements provides a good covariance matrix

Usage

```
validateCovariance(resVar, correlation, totalVar = NULL)
```

Arguments

resVar A vector of residual variances
correlation A correlation matrix
totalVar A vector of total variances

Value

Return TRUE if the covariance matrix is good

Author(s)

Sunthud Pornprasertmanit (University of Kansas; psunthud@ku.edu)

Examples

```
# No example
```

validateObject

Validate whether the drawn parameters are good.

Description

Validate whether the drawn parameters are good (providing an identified model).

Usage

```
validateObject(paramSet)
```

Arguments

paramSet A target set of parameters

Value

Return TRUE if the target parameters are good.

Author(s)

Sunthud Pornprasertmanit (University of Kansas; psunthud@ku.edu)

Examples

```
# No example
```

| | |
|--------------|--|
| validatePath | <i>Validate whether the regression coefficient (or loading) matrix is good</i> |
|--------------|--|

Description

Validate whether the regression coefficient (or loading) matrix is good

Usage

```
validatePath(path, var.iv, var.dv)
```

Arguments

| | |
|--------|---|
| path | A regression coefficient or loading matrix |
| var.iv | The variances of variables corresponding to the columns |
| var.dv | The variances of variables corresponding to the rows |

Value

Return TRUE if the target regression coefficient matrix is good.

Author(s)

Sunthud Pornprasertmanit (University of Kansas; psunthud@ku.edu)

Examples

```
# No example
```

| | |
|----------------|--|
| whichMonotonic | <i>Extract a part of a vector that is monotonically increasing or decreasing</i> |
|----------------|--|

Description

Extract a part of a vector that is monotonically increasing or decreasing. This function will go to the anchor value and extract the neighbor values that are monotonically increasing or decreasing.

Usage

```
whichMonotonic(vec, ord=NULL, anchor=NULL)
```

Arguments

| | |
|--------|---|
| vec | The target vector to be extracted |
| ord | The names of each element of the vector to be attached |
| anchor | The position of the element to be anchored. The default value is the middle position. |

Value

The monotonic part of a vector

Author(s)

Sunthud Pornprasertmanit (University of Kansas; <psunthud@ku.edu>)

Examples

```
# This is a private function.  
  
# whichMonotonic(c(3, 4, 1, 2, 3, 5, 2, 1))
```

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