

Package: lavaan (via r-universe)

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Title Latent Variable Analysis

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Description Fit a variety of latent variable models, including confirmatory factor analysis, structural equation modeling and latent growth curve models.

Depends R(>= 3.4)

Imports methods, stats4, stats, utils, graphics, MASS, mnormt, pbivnorm, numDeriv, quadprog

BugReports <https://github.com/yrosseel/lavaan/issues>

License GPL (>= 2)

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bootstrapLavaan	<i>Bootstrapping a Lavaan Model</i>
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Description

Bootstrap the LRT, or any other statistic (or vector of statistics) you can extract from a fitted lavaan object.

Usage

```
bootstrapLavaan(object, R = 1000L, type = "ordinary", verbose = FALSE,
  FUN = "coef", keep.idx = FALSE,
  parallel = c("no", "multicore", "snow"),
  ncpus = 1L, cl = NULL, iseed = NULL, h0.rmsea = NULL, ...)
```

```
bootstrapLRT(h0 = NULL, h1 = NULL, R = 1000L, type="bollen.stine",
  verbose = FALSE, return.LRT = FALSE, double.bootstrap = "no",
  double.bootstrap.R = 500L, double.bootstrap.alpha = 0.05,
  parallel = c("no", "multicore", "snow"),
  ncpus = 1L, cl = NULL, iseed = NULL)
```

Arguments

object	An object of class lavaan .
h0	An object of class lavaan . The restricted model.
h1	An object of class lavaan . The unrestricted model.
R	Integer. The number of bootstrap draws.
type	If "ordinary" or "nonparametric", the usual (naive) bootstrap method is used. If "bollen.stine", the data is first transformed such that the null hypothesis holds exactly in the resampling space. If "yuan", the data is first transformed by combining data and theory (model), such that the resampling space is closer to the population space. Note that both "bollen.stine" and "yuan" require the data to be continuous. They will not work with ordinal data. If "parametric", the parametric bootstrap approach is used; currently, this is only valid for continuous data following a multivariate normal distribution. See references for more details.
FUN	A function which when applied to the lavaan object returns a vector containing the statistic(s) of interest. The default is FUN="coef", returning the estimated values of the free parameters in the model.
...	Other named arguments for FUN which are passed unchanged each time it is called.
verbose	If TRUE, show information for each bootstrap draw.
keep.idx	If TRUE, store the indices of each bootstrap run (i.e., the observations that were used for this bootstrap run) as an attribute.

return.LRT	If TRUE, return the LRT values as an attribute to the pvalue.
parallel	The type of parallel operation to be used (if any). If missing, the default is "no".
ncpus	Integer: number of processes to be used in parallel operation. By default this is the number of cores (as detected by <code>parallel::detectCores()</code>) minus one.
cl	An optional parallel or snow cluster for use if <code>parallel = "snow"</code> . If not supplied, a cluster on the local machine is created for the duration of the <code>bootstrapLavaan</code> or <code>bootstrapLRT</code> call.
iseed	An integer to set the seed. Or NULL if no reproducible results are needed. This works for both serial (non-parallel) and parallel settings. Internally, <code>RNGkind()</code> is set to "L'Ecuyer-CMRG" if <code>parallel = "multicore"</code> . If <code>parallel = "snow"</code> (under windows), <code>parallel::clusterSetRNGStream()</code> is called which automatically switches to "L'Ecuyer-CMRG". When <code>iseed</code> is not NULL, <code>.Random.seed</code> (if it exists) in the global environment is left untouched.
h0.rmsea	Only used if <code>type="yuan"</code> . Allows one to do the Yuan bootstrap under the hypothesis that the population RMSEA equals a specified value.
double.bootstrap	If "standard" the genuine double bootstrap is used to compute an additional set of plug-in p-values for each bootstrap sample. If "FDB", the fast double bootstrap is used to compute second level LRT-values for each bootstrap sample. If "no", no double bootstrap is used. The default is set to "FDB".
double.bootstrap.R	Integer. The number of bootstrap draws to be use for the double bootstrap.
double.bootstrap.alpha	The significance level to compute the adjusted alpha based on the plugin p-values.

Details

The FUN function can return either a scalar or a numeric vector. This function can be an existing function (for example `coef`) or can be a custom defined function. For example:

```
myFUN <- function(x) {
  # require(lavaan)
  modelImpliedCov <- fitted(x)$cov
  vech(modelImpliedCov)
}
```

If `parallel="snow"`, it is imperative that the `require(lavaan)` is included in the custom function.

Value

For `bootstrapLavaan()`, the bootstrap distribution of the value(s) returned by FUN, when the object can be simplified to a vector. For `bootstrapLRT()`, a bootstrap *p* value, calculated as the proportion of bootstrap samples with a LRT statistic at least as large as the LRT statistic for the original data.

Author(s)

Yves Rosseel and Leonard Vanbrabant. Ed Merkle contributed Yuan's bootstrap. Improvements to Yuan's bootstrap were contributed by Hao Wu and Chuchu Cheng. The handling of iseed was contributed by Shu Fai Cheung.

References

Bollen, K. and Stine, R. (1992) Bootstrapping Goodness of Fit Measures in Structural Equation Models. *Sociological Methods and Research*, 21, 205–229.

Yuan, K.-H., Hayashi, K., & Yanagihara, H. (2007). A class of population covariance matrices in the bootstrap approach to covariance structure analysis. *Multivariate Behavioral Research*, 42, 261–281.

Examples

```
# fit the Holzinger and Swineford (1939) example
HS.model <- ' visual  =~ x1 + x2 + x3
             textual =~ x4 + x5 + x6
             speed   =~ x7 + x8 + x9 '

fit <- cfa(HS.model, data=HolzingerSwineford1939, se="none")

# get the test statistic for the original sample
T.orig <- fitMeasures(fit, "chisq")

# bootstrap to get bootstrap test statistics
# we only generate 10 bootstrap sample in this example; in practice
# you may wish to use a much higher number
T.boot <- bootstrapLavaan(fit, R=10, type="bollen.stine",
                        FUN=fitMeasures, fit.measures="chisq")

# compute a bootstrap based p-value
pvalue.boot <- length(which(T.boot > T.orig))/length(T.boot)
```

cfa

Fit Confirmatory Factor Analysis Models

Description

Fit a Confirmatory Factor Analysis (CFA) model.

Usage

```
cfa(model = NULL, data = NULL, ordered = NULL, sampling.weights = NULL,
     sample.cov = NULL, sample.mean = NULL, sample.th = NULL,
     sample.nobs = NULL, group = NULL, cluster = NULL,
     constraints = "", WLS.V = NULL, NACOV = NULL, ov.order = "model",
     ...)
```

Arguments

<code>model</code>	A description of the user-specified model. Typically, the model is described using the lavaan model syntax. See model.syntax for more information. Alternatively, a parameter table (eg. the output of the <code>lavaanify()</code> function) is also accepted.
<code>data</code>	An optional data frame containing the observed variables used in the model. If some variables are declared as ordered factors, lavaan will treat them as ordinal variables.
<code>ordered</code>	Character vector. Only used if the data is in a data.frame. Treat these variables as ordered (ordinal) variables, if they are endogenous in the model. Importantly, all other variables will be treated as numeric (unless they are declared as ordered in the data.frame.) Since 0.6-4, ordered can also be logical. If TRUE, all observed endogenous variables are treated as ordered (ordinal). If FALSE, all observed endogenous variables are considered to be numeric (again, unless they are declared as ordered in the data.frame.)
<code>sampling.weights</code>	A variable name in the data frame containing sampling weight information. Currently only available for non-clustered data. Depending on the <code>sampling.weights.normalization</code> option, these weights may be rescaled (or not) so that their sum equals the number of observations (total or per group).
<code>sample.cov</code>	Numeric matrix. A sample variance-covariance matrix. The rownames and/or colnames must contain the observed variable names. For a multiple group analysis, a list with a variance-covariance matrix for each group.
<code>sample.mean</code>	A sample mean vector. For a multiple group analysis, a list with a mean vector for each group.
<code>sample.th</code>	Vector of sample-based thresholds. For a multiple group analysis, a list with a vector of thresholds for each group.
<code>sample.nobs</code>	Number of observations if the full data frame is missing and only sample moments are given. For a multiple group analysis, a list or a vector with the number of observations for each group.
<code>group</code>	Character. A variable name in the data frame defining the groups in a multiple group analysis.
<code>cluster</code>	Character. A (single) variable name in the data frame defining the clusters in a two-level dataset.
<code>constraints</code>	Additional (in)equality constraints not yet included in the model syntax. See model.syntax for more information.
<code>WLS.V</code>	A user provided weight matrix to be used by estimator "WLS"; if the estimator is "DWLS", only the diagonal of this matrix will be used. For a multiple group analysis, a list with a weight matrix for each group. The elements of the weight matrix should be in the following order (if all data is continuous): first the means (if a meanstructure is involved), then the lower triangular elements of the covariance matrix including the diagonal, ordered column by column. In the categorical case: first the thresholds (including the means for continuous variables), then the slopes (if any), the variances of continuous variables (if any), and finally the lower triangular elements of the correlation/covariance matrix excluding the diagonal, ordered column by column.

NACOV	A user provided matrix containing the elements of (N times) the asymptotic variance-covariance matrix of the sample statistics. For a multiple group analysis, a list with an asymptotic variance-covariance matrix for each group. See the <code>WLS.V</code> argument for information about the order of the elements.
ov.order	Character. If "model" (the default), the order of the observed variable names (as reflected for example in the output of <code>lavNames()</code>) is determined by the model syntax. If "data", the order is determined by the data (either the full data.frame or the sample (co)variance matrix). If the <code>WLS.V</code> and/or <code>NACOV</code> matrices are provided, this argument is currently set to "data".
...	Many more additional options can be defined, using 'name = value'. See lavOptions for a complete list.

Details

The `cfa` function is a wrapper for the more general [lavaan](#) function, using the following default arguments: `int.ov.free = TRUE`, `int.lv.free = FALSE`, `auto.fix.first = TRUE` (unless `std.lv = TRUE`), `auto.fix.single = TRUE`, `auto.var = TRUE`, `auto.cov.lv.x = TRUE`, `auto.efa = TRUE`, `auto.th = TRUE`, `auto.delta = TRUE`, and `auto.cov.y = TRUE`.

Value

An object of class [lavaan](#), for which several methods are available, including a summary method.

References

Yves Rosseel (2012). `lavaan`: An R Package for Structural Equation Modeling. *Journal of Statistical Software*, 48(2), 1-36. doi:[10.18637/jss.v048.i02](https://doi.org/10.18637/jss.v048.i02)

See Also

[lavaan](#)

Examples

```
## The famous Holzinger and Swineford (1939) example
HS.model <- ' visual =~ x1 + x2 + x3
             textual =~ x4 + x5 + x6
             speed  =~ x7 + x8 + x9 '

fit <- cfa(HS.model, data = HolzingerSwineford1939)
summary(fit, fit.measures = TRUE)
```

`Demo.growth`*Demo dataset for a illustrating a linear growth model.*

Description

A toy dataset containing measures on 4 time points (t1,t2, t3 and t4), two predictors (x1 and x2) influencing the random intercept and slope, and a time-varying covariate (c1, c2, c3 and c4).

Usage

```
data(Demo.growth)
```

Format

A data frame of 400 observations of 10 variables.

t1 Measured value at time point 1

t2 Measured value at time point 2

t3 Measured value at time point 3

t4 Measured value at time point 4

x1 Predictor 1 influencing intercept and slope

x2 Predictor 2 influencing intercept and slope

c1 Time-varying covariate time point 1

c2 Time-varying covariate time point 2

c3 Time-varying covariate time point 3

c4 Time-varying covariate time point 4

See Also

[growth](#)

Examples

```
head(Demo.growth)
```

`Demo.twolevel`*Demo dataset for a illustrating a multilevel CFA.*

Description

A toy dataset containing measures on 6 items (y1-y6), 3 within-level covariates (x1-x3) and 2 between-level covariates (w1-w2). The data is clustered (200 clusters of size 5, 10, 15 and 20), and the cluster variable is “cluster”.

Usage

```
data(Demo.twolevel)
```

Format

A data frame of 2500 observations of 12 variables. clusters.

```
y1 item 1
y2 item 2
y3 item 3
y4 item 4
y5 item 5
y6 item 6
x1 within-level covariate 1
x2 within-level covariate 2
x3 within-level covariate 3
w1 between-level covariate 1
w2 between-level covariate 2
cluster cluster variable
```

Examples

```
head(Demo.twolevel)

model <- '
  level: 1
    fw =~ y1 + y2 + y3
    fw ~ x1 + x2 + x3
  level: 2
    fb =~ y1 + y2 + y3
    fb ~ w1 + w2
'

fit <- sem(model, data = Demo.twolevel, cluster = "cluster")
summary(fit)
```

Description

Fit one or more Exploratory Factor Analysis (EFA) model(s).

Usage

```
efa(data = NULL, nfactores = 1L, sample.cov = NULL, sample.nobs = NULL,
     rotation = "geomin", rotation.args = list(), ov.names = names(data),
     bounds = "pos.var", ..., output = "efa")
```

Arguments

<code>data</code>	A data frame containing the observed variables we need for the EFA. If only a subset of the observed variables is needed, use the <code>ov.names</code> argument.
<code>nfactores</code>	Integer or Integer vector. The desired number of factors to extract. Can be a single number, or a vector of numbers (e.g., <code>nfactores = 1:4.</code>), For each different number, a model is fitted.
<code>sample.cov</code>	Numeric matrix. A sample variance-covariance matrix. The rownames and/or colnames must contain the observed variable names. Unlike <code>sem</code> and <code>CFA</code> , the matrix may be a correlation matrix.
<code>sample.nobs</code>	Number of observations if the full data frame is missing and only the sample variance-covariance matrix is given.
<code>rotation</code>	Character. The rotation method to be used. Possible options are <code>varimax</code> , <code>quartimax</code> , <code>orthomax</code> , <code>oblimin</code> , <code>quartimin</code> , <code>geomin</code> , <code>promax</code> , <code>entropy</code> , <code>mccammon</code> , <code>infomax</code> , <code>tandem1</code> , <code>tandem2</code> , <code>oblimax</code> , <code>bentler</code> , <code>simplimax</code> , <code>target</code> , <code>pst</code> (=partially specified target), <code>cf</code> , <code>crawford-ferguson</code> , <code>cf-quartimax</code> , <code>cf-varimax</code> , <code>cf-equamax</code> , <code>cf-parsimax</code> , <code>cf-facparsim</code> , <code>biquartimin</code> , <code>bigeomin</code> . The latter two are for bifactor rotation only. The rotation algorithms (except <code>promax</code>) are similar to those from the <code>GPArotation</code> package, but have been reimplemented for better control. The <code>promax</code> method is taken from the <code>stats</code> package.
<code>rotation.args</code>	List. Options related to the rotation algorithm. The default options (and their alternatives) are <code>orthogonal = FALSE</code> , <code>row.weights = "default"</code> (or <code>"kaiser"</code> , <code>"cureton.mulaik"</code> or <code>"none"</code>), <code>std.ov = TRUE</code> , <code>algorithm = "gpa"</code> (or <code>"pairwise"</code>), <code>rstarts = 30</code> , <code>gpa.tol = 1e-05</code> , <code>tol = 1e-08</code> , <code>max.iter = 10000L</code> , <code>warn = FALSE</code> , <code>verbose = FALSE</code> , <code>reflect = TRUE</code> , <code>order.lv.by = "index"</code> (or <code>"sumofsquares"</code> or <code>"none"</code>). Other options are specific for a particular rotation criterion: <code>geomin.epsilon = 0.001</code> , <code>orthomax.gamma = 1</code> , <code>promax.kappa = 4</code> , <code>cf.gamma = 0</code> , and <code>oblimin.gamma = 0</code> .
<code>ov.names</code>	Character vector. The variables names that are needed for the EFA. Should be a subset of the variables names in the <code>data.frame</code> . By default, all the variables in the data are used.

bounds	Per default, bounds = "pos.var" forces all variances of both observed and latent variables to be strictly nonnegative. See the entry in lavOptions for more options.
...	Additional options to be passed to lavaan, using 'name = value'. See lavOptions for a complete list.
output	Character. If "efa" (the default), the output mimics the typical output of an EFA. If "lavaan", a lavaan object returned. The latter is only possible if nfactores contains a single (integer) number.

Details

The efa function is essentially a wrapper around the lavaan function. It generates the model syntax (for a given number of factors) and then calls lavaan() treating the factors as a single block that should be rotated. The function only supports a single group. Categorical data is handled as usual by first computing an appropriate (e.g., tetrachoric or polychoric) correlation matrix, which is then used as input for the EFA. There is also (limited) support for twolevel data. The same number of factors is then extracted at the within and the between level. The promax rotation method (taken from the stats package) is only provided for convenience. Because promax is a two-step algorithm (first varimax, then oblique rotation to get simple structure), it does not use the gpa or pairwise rotation algorithms, and as a result, no standard errors are provided.

Value

If output = "lavaan", an object of class [lavaan](#). If output = "efa", a list of class efaList for which a print(), summary() and fitMeasures() method are available. Because we added the (standardized) loadings as an extra element, the loadings function (which is not a generic function) from the stats package will also work on efaList objects.

See Also

[summary.efaList](#) for a summary method if the output is of class efaList.

Examples

```
## The famous Holzinger and Swineford (1939) example
fit <- efa(data = HolzingerSwineford1939,
          ov.names = paste("x", 1:9, sep = ""),
          nfactors = 1:3,
          rotation = "geomin",
          rotation.args = list(geomin.epsilon = 0.01, rstarts = 1))
summary(fit, nd = 3L, cutoff = 0.2, dot.cutoff = 0.05)
fitMeasures(fit, fit.measures = "all")
```

 estfun

Extract Empirical Estimating Functions

Description

A function for extracting the empirical estimating functions of a fitted lavaan model. This is the derivative of the objective function with respect to the parameter vector, evaluated at the observed (case-wise) data. In other words, this function returns the case-wise scores, evaluated at the fitted model parameters.

Usage

```
estfun.lavaan(object, scaling = FALSE, ignore.constraints = FALSE,
              remove.duplicated = TRUE, remove.empty.cases = TRUE)
lavScores(object, scaling = FALSE, ignore.constraints = FALSE,
           remove.duplicated = TRUE, remove.empty.cases = TRUE)
```

Arguments

<code>object</code>	An object of class <code>lavaan</code> .
<code>scaling</code>	Only used for the ML estimator. If TRUE, the scores are scaled to reflect the specific objective function used by lavaan. If FALSE (the default), the objective function is the loglikelihood function assuming multivariate normality.
<code>ignore.constraints</code>	Logical. If TRUE, the scores do not reflect the (equality or inequality) constraints. If FALSE, the scores are computed by taking the unconstrained scores, and adding the term $t(R) \lambda$, where λ are the (case-wise) Lagrange Multipliers, and R is the Jacobian of the constraint function. Only in the latter case will the sum of the columns be (almost) equal to zero.
<code>remove.duplicated</code>	If TRUE, and all the equality constraints have a simple form (eg. $a == b$), the unconstrained scores are post-multiplied with a transformation matrix in order to remove the duplicated parameters.
<code>remove.empty.cases</code>	If TRUE, empty cases with only missing values will be removed from the output.

Value

A $n \times k$ matrix corresponding to n observations and k parameters.

Author(s)

Ed Merkle for the ML case; the `remove.duplicated`, `ignore.constraints` and `remove.empty.cases` arguments were added by Yves Rosseel; Franz Classe for the WLS case.

Examples

```
## The famous Holzinger and Swineford (1939) example
HS.model <- ' visual  =~ x1 + x2 + x3
             textual =~ x4 + x5 + x6
             speed   =~ x7 + x8 + x9 '

fit <- cfa(HS.model, data = HolzingerSwineford1939)
head(lavScores(fit))
```

FacialBurns

Dataset for illustrating the InformativeTesting function.

Description

A dataset from the Dutch burn center (<http://www.adbc.nl>). The data were used to examine psychosocial functioning in patients with facial burn wounds. Psychosocial functioning was measured by Anxiety and depression symptoms (HADS), and self-esteem (Rosenberg's self-esteem scale).

Usage

```
data(FacialBurns)
```

Format

A data frame of 77 observations of 6 variables.

Selfesteem Rosenberg's self-esteem scale

HADS Anxiety and depression scale

Age Age measured in years, control variable

TBSA Total Burned Surface Area

RUM Rumination, control variable

Sex Gender, grouping variable

Examples

```
head(FacialBurns)
```

Description

This function computes a variety of fit measures to assess the global fit of a latent variable model.

Usage

```
fitMeasures(object, fit.measures = "all",
            baseline.model = NULL, h1.model = NULL,
            fm.args = list(standard.test = "default",
                          scaled.test = "default",
                          rmsea.ci.level = 0.90,
                          rmsea.close.h0 = 0.05,
                          rmsea.notclose.h0 = 0.08,
                          robust = TRUE,
                          cat.check.pd = TRUE),
            output = "vector", ...)
fitmeasures(object, fit.measures = "all",
            baseline.model = NULL, h1.model = NULL,
            fm.args = list(standard.test = "default",
                          scaled.test = "default",
                          rmsea.ci.level = 0.90,
                          rmsea.close.h0 = 0.05,
                          rmsea.notclose.h0 = 0.08,
                          robust = TRUE,
                          cat.check.pd = TRUE),
            output = "vector", ...)
```

Arguments

- | | |
|----------------|---|
| object | An object of class lavaan . |
| fit.measures | If "all", all fit measures available will be returned. If only a single or a few fit measures are specified by name, only those are computed and returned. |
| baseline.model | If not NULL, an object of class lavaan , representing a user-specified baseline model. If a baseline model is provided, all fit indices relying on a baseline model (eg. CFI or TLI) will use the test statistics from this user-specified baseline model, instead of the default baseline model. |
| h1.model | If not NULL, an object of class lavaan , representing a user-specified alternative to the default unrestricted model. If h1.model is provided, all fit indices calculated from chi-squared will use the chi-squared <i>difference</i> test statistics from lavTestLRT , which compare the user-provided h1.model to object. |
| fm.args | List. Additional options for certain fit measures. The standard.test element determines the main test statistic (chi-square value) that will be used to compute all the fit measures that depend on this test statistic. Usually this is "standard". |

The `scaled.test` element determines which scaling method is to be used for the scaled fit measures (in case multiple scaling methods were requested). The `rmsea.ci.level` element determines the level of the confidence interval for the `rmsea` value. The `rmsea.close.h0` element is the `rmsea` value that is used under the null hypothesis that $rmsea \leq rmsea.close.h0$. The `rmsea.notclose.h0` element is the `rmsea` value that is used under the null hypothesis that $rmsea >= rmsea.notclose.h0$. The `robust` element can be set to `FALSE` to avoid computing the so-called robust `rmsea/cfi` measures (for example if the computations take too long). The `cat.check.pd` element is only used when data is categorical. If `TRUE`, robust values for RMSEA and CFI are only computed if the input correlation matrix is positive-definite (for all groups).

output	Character. If "vector" (the default), display the output as a named (lavaan-formatted) vector. If "matrix", display the output as a 1-column matrix. If "text", display the output using subsections and verbose descriptions. The latter is used in the summary output, and does not print the chi-square test by default. In addition, <code>fit.measures</code> should contain the main ingredient (for example "rmsea") if related fit measures are requested (for example "rmsea.ci.lower"). Otherwise, nothing will be printed in that section. See the examples how to add the chi-square test in the text output.
...	Further arguments passed to or from other methods. Not currently used for lavaan objects.

Value

A named numeric vector of fit measures.

Examples

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

```
fit <- cfa(HS.model, data = HolzingerSwineford1939)
fitMeasures(fit)
fitMeasures(fit, "cfi")
fitMeasures(fit, c("chisq", "df", "pvalue", "cfi", "rmsea"))
fitMeasures(fit, c("chisq", "df", "pvalue", "cfi", "rmsea"),
             output = "matrix")
fitMeasures(fit, c("chisq", "df", "pvalue", "cfi", "rmsea"),
             output = "text")

## fit a more restricted model
fit0 <- cfa(HS.model, data = HolzingerSwineford1939, orthogonal = TRUE)
## Calculate RMSEA_D (Savalei et al., 2023)
## See https://psycnet.apa.org/doi/10.1037/met0000537
fitMeasures(fit0, "rmsea", h1.model = fit)
```

getCov

*Utility Functions For Covariance Matrices***Description**

Convenience functions to deal with covariance and correlation matrices.

Usage

```
getCov(x, lower = TRUE, diagonal = TRUE, sds = NULL,
       names = paste("V", 1:nvar, sep=""))
char2num(s)
cor2cov(R, sds, names = NULL)
```

Arguments

x	The elements of the covariance matrix. Either inside a character string or as a numeric vector. In the former case, the function <code>char2num</code> is used to convert the numbers (inside the character string) to numeric values.
lower	Logical. If <code>TRUE</code> , the numeric values in <code>x</code> are the lower-triangular elements of the (symmetric) covariance matrix only. If <code>FALSE</code> , <code>x</code> contains the upper triangular elements only. Note we always assumed the elements are provided row-wise!
diagonal	Logical. If <code>TRUE</code> , the numeric values in <code>x</code> include the diagonal elements. If <code>FALSE</code> , a unit diagonal is assumed.
sds	A numeric vector containing the standard deviations to be used to scale the elements in <code>x</code> or the correlation matrix <code>R</code> into a covariance matrix.
names	The variable names of the observed variables.
s	Character string containing numeric values; comma's and semi-colons are ignored.
R	A correlation matrix, to be scaled into a covariance matrix.

Details

The `getCov` function is typically used to input the lower (or upper) triangular elements of a (symmetric) covariance matrix. In many examples found in handbooks, only those elements are shown. However, `lavaan` needs a full matrix to proceed.

The `cor2cov` function is the inverse of the `cov2cor` function, and scales a correlation matrix into a covariance matrix given the standard deviations of the variables. Optionally, variable names can be given.

Examples

```
# The classic Wheaton et. al. (1977) model
# panel data on the stability of alienation
lower <- '
```



```

11.834,
 6.947,   9.364,
 6.819,   5.091,  12.532,
 4.783,   5.028,   7.495,   9.986,
-3.839,  -3.889,  -3.841,  -3.625,   9.610,
-21.899, -18.831, -21.748, -18.775,  35.522,  450.288 '

# convert to a full symmetric covariance matrix with names
wheaton.cov <- getCov(lower, names=c("anomia67", "powerless67", "anomia71",
                                     "powerless71", "education", "sei"))

# the model
wheaton.model <- '
# measurement model
  ses      =~ education + sei
  alien67 =~ anomia67 + powerless67
  alien71 =~ anomia71 + powerless71

# equations
  alien71 ~ alien67 + ses
  alien67 ~ ses

# correlated residuals
  anomia67 ~~ anomia71
  powerless67 ~~ powerless71
,

# fitting the model
fit <- sem(wheaton.model, sample.cov=wheaton.cov, sample.nobs=932)

# showing the results
summary(fit, standardized=TRUE)

```

growth

Fit Growth Curve Models

Description

Fit a Growth Curve model. Only useful if all the latent variables in the model are growth factors. For more complex models, it may be better to use the [lavaan](#) function.

Usage

```

growth(model = NULL, data = NULL, ordered = NULL, sampling.weights = NULL,
        sample.cov = NULL, sample.mean = NULL, sample.th = NULL,
        sample.nobs = NULL, group = NULL, cluster = NULL,
        constraints = "", WLS.V = NULL, NACOV = NULL, ov.order = "model",
        ...)

```

Arguments

<code>model</code>	A description of the user-specified model. Typically, the model is described using the lavaan model syntax. See model.syntax for more information. Alternatively, a parameter table (eg. the output of the <code>lavaanify()</code> function) is also accepted.
<code>data</code>	An optional data frame containing the observed variables used in the model. If some variables are declared as ordered factors, lavaan will treat them as ordinal variables.
<code>ordered</code>	Character vector. Only used if the data is in a data.frame. Treat these variables as ordered (ordinal) variables, if they are endogenous in the model. Importantly, all other variables will be treated as numeric (unless they are declared as ordered in the data.frame.) Since 0.6-4, ordered can also be logical. If TRUE, all observed endogenous variables are treated as ordered (ordinal). If FALSE, all observed endogenous variables are considered to be numeric (again, unless they are declared as ordered in the data.frame.)
<code>sampling.weights</code>	A variable name in the data frame containing sampling weight information. Currently only available for non-clustered data. Depending on the <code>sampling.weights.normalization</code> option, these weights may be rescaled (or not) so that their sum equals the number of observations (total or per group).
<code>sample.cov</code>	Numeric matrix. A sample variance-covariance matrix. The rownames and/or colnames must contain the observed variable names. For a multiple group analysis, a list with a variance-covariance matrix for each group.
<code>sample.mean</code>	A sample mean vector. For a multiple group analysis, a list with a mean vector for each group.
<code>sample.th</code>	Vector of sample-based thresholds. For a multiple group analysis, a list with a vector of thresholds for each group.
<code>sample.nobs</code>	Number of observations if the full data frame is missing and only sample moments are given. For a multiple group analysis, a list or a vector with the number of observations for each group.
<code>group</code>	Character. A variable name in the data frame defining the groups in a multiple group analysis.
<code>cluster</code>	Character. A (single) variable name in the data frame defining the clusters in a two-level dataset.
<code>constraints</code>	Additional (in)equality constraints not yet included in the model syntax. See model.syntax for more information.
<code>WLS.V</code>	A user provided weight matrix to be used by estimator "WLS"; if the estimator is "DWLS", only the diagonal of this matrix will be used. For a multiple group analysis, a list with a weight matrix for each group. The elements of the weight matrix should be in the following order (if all data is continuous): first the means (if a meanstructure is involved), then the lower triangular elements of the covariance matrix including the diagonal, ordered column by column. In the categorical case: first the thresholds (including the means for continuous variables), then the slopes (if any), the variances of continuous variables (if any), and finally the lower triangular elements of the correlation/covariance matrix excluding the diagonal, ordered column by column.

NACOV	A user provided matrix containing the elements of (N times) the asymptotic variance-covariance matrix of the sample statistics. For a multiple group analysis, a list with an asymptotic variance-covariance matrix for each group. See the WLS.V argument for information about the order of the elements.
ov.order	Character. If "model" (the default), the order of the observed variable names (as reflected for example in the output of <code>lavNames()</code>) is determined by the model syntax. If "data", the order is determined by the data (either the full data.frame or the sample (co)variance matrix). If the WLS.V and/or NACOV matrices are provided, this argument is currently set to "data".
...	Many more additional options can be defined, using 'name = value'. See lavOptions for a complete list.

Details

The growth function is a wrapper for the more general [lavaan](#) function, using the following default arguments: `meanstructure = TRUE`, `int.ov.free = FALSE`, `int.lv.free = TRUE`, `auto.fix.first = TRUE` (unless `std.lv = TRUE`), `auto.fix.single = TRUE`, `auto.var = TRUE`, `auto.cov.lv.x = TRUE`, `auto.efa = TRUE`, `auto.th = TRUE`, `auto.delta = TRUE`, and `auto.cov.y = TRUE`.

Value

An object of class [lavaan](#), for which several methods are available, including a summary method.

References

Yves Rosseel (2012). [lavaan: An R Package for Structural Equation Modeling](#). *Journal of Statistical Software*, 48(2), 1-36. doi:[10.18637/jss.v048.i02](https://doi.org/10.18637/jss.v048.i02)

See Also

[lavaan](#)

Examples

```
## linear growth model with a time-varying covariate
model.syntax <- '
  # intercept and slope with fixed coefficients
  i =~ 1*t1 + 1*t2 + 1*t3 + 1*t4
  s =~ 0*t1 + 1*t2 + 2*t3 + 3*t4

  # regressions
  i ~ x1 + x2
  s ~ x1 + x2

  # time-varying covariates
  t1 ~ c1
  t2 ~ c2
  t3 ~ c3
  t4 ~ c4
'
```

```
fit <- growth(model.syntax, data = Demo.growth)
summary(fit)
```

HolzingerSwineford1939

Holzinger and Swineford Dataset (9 Variables)

Description

The classic Holzinger and Swineford (1939) dataset consists of mental ability test scores of seventh- and eighth-grade children from two different schools (Pasteur and Grant-White). In the original dataset (available in the MBESS package), there are scores for 26 tests. However, a smaller subset with 9 variables is more widely used in the literature (for example in Joreskog's 1969 paper, which also uses the 145 subjects from the Grant-White school only).

Usage

```
data(HolzingerSwineford1939)
```

Format

A data frame with 301 observations of 15 variables.

id Identifier
sex Gender
ageyr Age, year part
agemo Age, month part
school School (Pasteur or Grant-White)
grade Grade
x1 Visual perception
x2 Cubes
x3 Lozenges
x4 Paragraph comprehension
x5 Sentence completion
x6 Word meaning
x7 Speeded addition
x8 Speeded counting of dots
x9 Speeded discrimination straight and curved capitals

Source

This dataset was originally retrieved from <http://web.missouri.edu/~kolenikovs/stata/hs-cfa.dta> (link no longer active) and converted to an R dataset.

References

- Holzinger, K., and Swineford, F. (1939). A study in factor analysis: The stability of a bifactor solution. Supplementary Educational Monograph, no. 48. Chicago: University of Chicago Press.
- Joreskog, K. G. (1969). A general approach to confirmatory maximum likelihood factor analysis. *Psychometrika*, 34, 183-202.

See Also

[cfa](#)

Examples

```
head(HolzingerSwineford1939)
```

InformativeTesting *Testing order/inequality Constrained Hypotheses in SEM*

Description

Testing order/inequality constrained Hypotheses in SEM

Usage

```
InformativeTesting(model = NULL, data, constraints = NULL,
                   R = 1000L, type = "bollen.stine",
                   return.LRT = TRUE,
                   double.bootstrap = "standard",
                   double.bootstrap.R = 249L,
                   double.bootstrap.alpha = 0.05,
                   parallel = c("no", "multicore", "snow"),
                   ncpus = 1L, cl = NULL, verbose = FALSE, ...)
```

Arguments

<code>model</code>	Model syntax specifying the model. See model.syntax for more information.
<code>data</code>	The data frame containing the observed variables being used to fit the model.
<code>constraints</code>	The imposed inequality constraints on the model.
<code>R</code>	Integer; number of bootstrap draws. The default value is set to 1000.
<code>type</code>	If "parametric", the parametric bootstrap is used. If "bollen.stine", the semi-nonparametric Bollen-Stine bootstrap is used. The default is set to "bollen.stine".
<code>return.LRT</code>	Logical; if TRUE, the function returns bootstrapped LRT-values.
<code>double.bootstrap</code>	If "standard" (default) the genuine double bootstrap is used to compute an additional set of plug-in p-values for each bootstrap sample. If "no", no double bootstrap is used. If "FDB", the fast double bootstrap is used to compute second level LRT-values for each bootstrap sample. Note that the "FDB" is experimental and should not be used by inexperienced users.

<code>double.bootstrap.R</code>	Integer; number of double bootstrap draws. The default value is set to 249.
<code>double.bootstrap.alpha</code>	The significance level to compute the adjusted alpha based on the plugin p-values. Only used if <code>double.bootstrap = "standard"</code> . The default value is set to 0.05.
<code>parallel</code>	The type of parallel operation to be used (if any). If missing, the default is set "no".
<code>ncpus</code>	Integer: number of processes to be used in parallel operation: typically one would chose this to the number of available CPUs.
<code>cl</code>	An optional parallel or snow cluster for use if <code>parallel = "snow"</code> . If not supplied, a cluster on the local machine is created for the duration of the <code>InformativeTesting</code> call.
<code>verbose</code>	Logical; if TRUE, information is shown at each bootstrap draw.
<code>...</code>	Other named arguments from the lavaan package which are passed to the function. For example "group" in a multiple group model.

Details

The following hypothesis tests are available:

- Type A: Test H0: all restriktions with equalities ("=") active against HA: at least one inequality restriktion (">") strictly true.
- Type B: Test H0: all restriktions with inequalities ">" (including some equalities "=") active against HA: at least one restriktion false (some equality restriktions may be maintained).

Value

An object of class `InformativeTesting` for which a `print` and a `plot` method is available.

Author(s)

Leonard Vanbrabant <lgf.vanbrabant@gmail.com>

References

- Van de Schoot, R., Hoijtink, H., & Dekovic, M. (2010). Testing inequality constrained hypotheses in SEM models. *Structural Equation Modeling*, **17**, 443-463.
- Van de Schoot, R., Strohmeier, D. (2011). Testing informative hypotheses in SEM increases power: An illustration contrasting classical. *International Journal of Behavioral Development*, **35**, 180-190.
- Silvapulle, M.J. and Sen, P.K. (2005). *Constrained Statistical Inference*. Wiley, New York.

Examples

```

## Not run:
#####
### real data example ###
#####
# Multiple group path model for facial burns example.

# model syntax with starting values.
burns.model <- 'Selfesteem ~ Age + c(m1, f1)*TBSA + HADS +
               start(-.10, -.20)*TBSA
               HADS ~ Age + c(m2, f2)*TBSA + RUM +
               start(.10, .20)*TBSA '

# constraints syntax
burns.constraints <- 'f2 > 0 ; m1 < 0
                    m2 > 0 ; f1 < 0
                    f2 > m2 ; f1 < m1'

# we only generate 2 bootstrap samples in this example; in practice
# you may wish to use a much higher number.
# the double bootstrap was switched off; in practice you probably
# want to set it to "standard".
example1 <- InformativeTesting(model = burns.model, data = FacialBurns,
                              R = 2, constraints = burns.constraints,
                              double.bootstrap = "no", group = "Sex")

example1

#####
### artificial example ###
#####
# Simple ANOVA model with 3 groups (N = 20 per group)
set.seed(1234)
Y <- cbind(c(rnorm(20,0,1), rnorm(20,0.5,1), rnorm(20,1,1)))
grp <- c(rep("1", 20), rep("2", 20), rep("3", 20))
Data <- data.frame(Y, grp)

#create model matrix
fit.lm <- lm(Y ~ grp, data = Data)
mfit <- fit.lm$model
mm <- model.matrix(mfit)

Y <- model.response(mfit)
X <- data.frame(mm[,2:3])
names(X) <- c("d1", "d2")
Data.new <- data.frame(Y, X)

# model
model <- 'Y ~ 1 + a1*d1 + a2*d2'

# fit without constraints

```

```

fit <- sem(model, data = Data.new)

# constraints syntax: mu1 < mu2 < mu3
constraints <- ' a1 > 0
                a1 < a2 '

# we only generate 10 bootstrap samples in this example; in practice
# you may wish to use a much higher number, say > 1000. The double
# bootstrap is not necessary in case of an univariate ANOVA model.
example2 <- InformativeTesting(model = model, data = Data.new,
                               start = parTable(fit),
                               R = 10L, double.bootstrap = "no",
                               constraints = constraints)

example2

## End(Not run)

```

InformativeTesting methods

Methods for output InformativeTesting()

Description

The print function shows the results of hypothesis tests Type A and Type B. The plot function plots the distributions of bootstrapped LRT values and plug-in p-values.

Usage

```

## S3 method for class 'InformativeTesting'
print(x, digits = max(3, getOption("digits") - 3), ...)

## S3 method for class 'InformativeTesting'
plot(x, ..., type = c("lr", "ppv"),
     main = "main", xlab = "xlabel", ylab = "Frequency", freq = TRUE,
     breaks = 15, cex.main = 1, cex.lab = 1, cex.axis = 1,
     col = "grey", border = par("fg"), vline = TRUE,
     vline.col = c("red", "blue"), lty = c(1,2), lwd = 1,
     legend = TRUE, bty = "o", cex.legend = 1, loc.legend = "topright")

```

Arguments

x	object of class "InformativeTesting".
digits	the number of significant digits to use when printing.
...	Currently not used.
type	If "lr", a distribution of the first-level bootstrapped LR values is plotted. If "ppv" a distribution of the bootstrapped plug-in p-values is plotted.
main	The main title(s) for the plot(s).

<code>xlab</code>	A label for the x axis, default depends on input type.
<code>ylab</code>	A label for the y axis.
<code>freq</code>	Logical; if TRUE, the histogram graphic is a representation of frequencies, the counts component of the result; if FALSE, probability densities, component density, are plotted (so that the histogram has a total area of one). The default is set to TRUE.
<code>breaks</code>	see hist
<code>cex.main</code>	The magnification to be used for main titles relative to the current setting of <code>cex</code> .
<code>cex.lab</code>	The magnification to be used for x and y labels relative to the current setting of <code>cex</code> .
<code>cex.axis</code>	The magnification to be used for axis annotation relative to the current setting of <code>cex</code> .
<code>col</code>	A colour to be used to fill the bars. The default of NULL yields unfilled bars.
<code>border</code>	Color for rectangle border(s). The default means <code>par("fg")</code> .
<code>vline</code>	Logical; if TRUE a vertical line is drawn at the observed LRT value. If <code>double.bootstrap = "FDB"</code> a vertical line is drawn at the 1-p* quantile of the second-level LRT values, where p* is the first-level bootstrapped p-value
<code>vline.col</code>	Color(s) for the <code>vline.LRT</code> .
<code>lty</code>	The line type. Line types can either be specified as an integer (0=blank, 1=solid (default), 2=dashed, 3=dotted, 4=dotdash, 5=longdash, 6=twodash) or as one of the character strings "blank", "solid", "dashed", "dotted", "dotdash", "longdash", or "twodash", where "blank" uses 'invisible lines' (i.e., does not draw them).
<code>lwd</code>	The line width, a positive number, defaulting to 1.
<code>legend</code>	Logical; if TRUE a legend is added to the plot.
<code>bty</code>	A character string which determined the type of box which is drawn about plots. If <code>bty</code> is one of "o" (the default), "l", "7", "c", "u", or "]" the resulting box resembles the corresponding upper case letter. A value of "n" suppresses the box.
<code>cex.legend</code>	A numerical value giving the amount by which the legend text and symbols should be magnified relative to the default. This starts as 1 when a device is opened, and is reset when the layout is changed.
<code>loc.legend</code>	The location of the legend, specified by a single keyword from the list "bottomright", "bottom", "bottomleft", "left", "topleft", "top", "topright", "right" and "center".

Author(s)

Leonard Vanbrabant <lgf.vanbrabant@gmail.com>

Examples

```
## Not run:
#####
### real data example ###
```

```
#####
# Multiple group path model for facial burns example.

# model syntax with starting values.
burns.model <- 'Selfesteem ~ Age + c(m1, f1)*TBSA + HADS +
               start(-.10, -.20)*TBSA
               HADS ~ Age + c(m2, f2)*TBSA + RUM +
               start(.10, .20)*TBSA '
```

```
# constraints syntax
burns.constraints <- 'f2 > 0 ; m1 < 0
                    m2 > 0 ; f1 < 0
                    f2 > m2 ; f1 < m1'
```

```
# we only generate 2 bootstrap samples in this example; in practice
# you may wish to use a much higher number.
# the double bootstrap was switched off; in practice you probably
# want to set it to "standard".
example1 <- InformativeTesting(model = burns.model, data = FacialBurns,
                              R = 2, constraints = burns.constraints,
                              double.bootstrap = "no", group = "Sex")

example1
plot(example1)

#####
### artificial example ###
#####
# Simple ANOVA model with 3 groups (N = 20 per group)
set.seed(1234)
Y <- cbind(c(rnorm(20,0,1), rnorm(20,0.5,1), rnorm(20,1,1)))
grp <- c(rep("1", 20), rep("2", 20), rep("3", 20))
Data <- data.frame(Y, grp)

#create model matrix
fit.lm <- lm(Y ~ grp, data = Data)
mfit <- fit.lm$model
mm <- model.matrix(mfit)

Y <- model.response(mfit)
X <- data.frame(mm[,2:3])
names(X) <- c("d1", "d2")
Data.new <- data.frame(Y, X)

# model
model <- 'Y ~ 1 + a1*d1 + a2*d2'

# fit without constraints
fit <- sem(model, data = Data.new)

# constraints syntax: mu1 < mu2 < mu3
constraints <- ' a1 > 0
               a1 < a2 '
```

```
# we only generate 10 bootstrap samples in this example; in practice
# you may wish to use a much higher number, say > 1000. The double
# bootstrap is not necessary in case of an univariate ANOVA model.
example2 <- InformativeTesting(model = model, data = Data.new,
                              start = parTable(fit),
                              R = 10L, double.bootstrap = "no",
                              constraints = constraints)

example2
# plot(example2)

## End(Not run)
```

`inspectSampleCov`*Observed Variable Correlation Matrix from a Model and Data*

Description

The lavaan model syntax describes a latent variable model. Often, the user wants to see the covariance matrix generated by their model for diagnostic purposes. However, their data may have far more columns of information than what is contained in their model.

Usage

```
inspectSampleCov(model, data, ...)
```

Arguments

<code>model</code>	The model that will be fit by lavaan.
<code>data</code>	The data frame being used to fit the model.
<code>...</code>	Other arguments to sem for how to deal with multiple groups, missing values, etc.

Details

One must supply both a model, coded with proper [model.syntax](#) and a data frame from which a covariance matrix will be calculated. This function essentially calls [sem](#), but doesn't fit the model, then uses [lavInspect](#) to get the sample covariance matrix and meanstructure.

See also

[sem](#), [lavInspect](#)

Author(s)

Jarrett Byrnes

lavaan

*Fit a Latent Variable Model***Description**

Fit a latent variable model.

Usage

```
lavaan(model = NULL, data = NULL, ordered = NULL,
        sampling.weights = NULL,
        sample.cov = NULL, sample.mean = NULL, sample.th = NULL,
        sample.nobs = NULL,
        group = NULL, cluster = NULL, constraints = "",
        WLS.V = NULL, NACOV = NULL, ov.order = "model",
        slotOptions = NULL, slotParTable = NULL, slotSampleStats = NULL,
        slotData = NULL, slotModel = NULL, slotCache = NULL,
        sloth1 = NULL,
        ...)
```

Arguments

model	A description of the user-specified model. Typically, the model is described using the lavaan model syntax. See model.syntax for more information. Alternatively, a parameter table (eg. the output of the <code>lavaanify()</code> function) is also accepted.
data	An optional data frame containing the observed variables used in the model. If some variables are declared as ordered factors, lavaan will treat them as ordinal variables.
ordered	Character vector. Only used if the data is in a data.frame. Treat these variables as ordered (ordinal) variables, if they are endogenous in the model. Importantly, all other variables will be treated as numeric (unless they are declared as ordered in the data.frame.) Since 0.6-4, ordered can also be logical. If TRUE, all observed endogenous variables are treated as ordered (ordinal). If FALSE, all observed endogenous variables are considered to be numeric (again, unless they are declared as ordered in the data.frame.)
sampling.weights	A variable name in the data frame containing sampling weight information. Currently only available for non-clustered data. Depending on the <code>sampling.weights.normalization</code> option, these weights may be rescaled (or not) so that their sum equals the number of observations (total or per group).
sample.cov	Numeric matrix. A sample variance-covariance matrix. The rownames and/or colnames must contain the observed variable names. For a multiple group analysis, a list with a variance-covariance matrix for each group.
sample.mean	A sample mean vector. For a multiple group analysis, a list with a mean vector for each group.

<code>sample.th</code>	Vector of sample-based thresholds. For a multiple group analysis, a list with a vector of thresholds for each group.
<code>sample.nobs</code>	Number of observations if the full data frame is missing and only sample moments are given. For a multiple group analysis, a list or a vector with the number of observations for each group.
<code>group</code>	Character. A variable name in the data frame defining the groups in a multiple group analysis.
<code>cluster</code>	Character. A (single) variable name in the data frame defining the clusters in a two-level dataset.
<code>constraints</code>	Additional (in)equality constraints not yet included in the model syntax. See model.syntax for more information.
<code>WLS.V</code>	A user provided weight matrix to be used by estimator "WLS"; if the estimator is "DWLS", only the diagonal of this matrix will be used. For a multiple group analysis, a list with a weight matrix for each group. The elements of the weight matrix should be in the following order (if all data is continuous): first the means (if a meanstructure is involved), then the lower triangular elements of the covariance matrix including the diagonal, ordered column by column. In the categorical case: first the thresholds (including the means for continuous variables), then the slopes (if any), the variances of continuous variables (if any), and finally the lower triangular elements of the correlation/covariance matrix excluding the diagonal, ordered column by column.
<code>NACOV</code>	A user provided matrix containing the elements of (N times) the asymptotic variance-covariance matrix of the sample statistics. For a multiple group analysis, a list with an asymptotic variance-covariance matrix for each group. See the <code>WLS.V</code> argument for information about the order of the elements.
<code>ov.order</code>	Character. If "model" (the default), the order of the observed variable names (as reflected for example in the output of <code>lavNames()</code>) is determined by the model syntax. If "data", the order is determined by the data (either the full data.frame or the sample (co)variance matrix). If the <code>WLS.V</code> and/or <code>NACOV</code> matrices are provided, this argument is currently set to "data".
<code>slotOptions</code>	Options slot from a fitted lavaan object. If provided, no new Options slot will be created by this call.
<code>slotParTable</code>	ParTable slot from a fitted lavaan object. If provided, no new ParTable slot will be created by this call.
<code>slotSampleStats</code>	SampleStats slot from a fitted lavaan object. If provided, no new SampleStats slot will be created by this call.
<code>slotData</code>	Data slot from a fitted lavaan object. If provided, no new Data slot will be created by this call.
<code>slotModel</code>	Model slot from a fitted lavaan object. If provided, no new Model slot will be created by this call.
<code>slotCache</code>	Cache slot from a fitted lavaan object. If provided, no new Cache slot will be created by this call.
<code>sloth1</code>	h1 slot from a fitted lavaan object. If provided, no new h1 slot will be created by this call.

... Many more additional options can be defined, using 'name = value'. See [lavOptions](#) for a complete list.

Value

An object of class [lavaan](#), for which several methods are available, including a summary method.

References

Yves Rosseel (2012). lavaan: An R Package for Structural Equation Modeling. Journal of Statistical Software, 48(2), 1-36. doi:10.18637/jss.v048.i02

See Also

[cfa](#), [sem](#), [growth](#)

Examples

```
# The Holzinger and Swineford (1939) example
HS.model <- ' visual =~ x1 + x2 + x3
             textual =~ x4 + x5 + x6
             speed  =~ x7 + x8 + x9 '

fit <- lavaan(HS.model, data=HolzingerSwineford1939,
              auto.var=TRUE, auto.fix.first=TRUE,
              auto.cov.lv.x=TRUE)
summary(fit, fit.measures=TRUE)
```

lavaan-class

Class For Representing A (Fitted) Latent Variable Model

Description

The lavaan class represents a (fitted) latent variable model. It contains a description of the model as specified by the user, a summary of the data, an internal matrix representation, and if the model was fitted, the fitting results.

Objects from the Class

Objects can be created via the [cfa](#), [sem](#), [growth](#) or [lavaan](#) functions.

Slots

version: The lavaan package version used to create this objects

call: The function call as returned by `match.call()`.

timing: The elapsed time (user+system) for various parts of the program as a list, including the total time.

Options: Named list of options that were provided by the user, or filled-in automatically.

ParTable: Named list describing the model parameters. Can be coerced to a data.frame. In the documentation, this is called the ‘parameter table’.

pta: Named list containing parameter table attributes.

Data: Object of internal class "Data": information about the data.

SampleStats: Object of internal class "SampleStats": sample statistics

Model: Object of internal class "Model": the internal (matrix) representation of the model

Cache: List using objects that we try to compute only once, and reuse many times.

Fit: Object of internal class "Fit": the results of fitting the model. No longer used.

boot: List. Results and information about the bootstrap.

optim: List. Information about the optimization.

loglik: List. Information about the loglikelihood of the model (if maximum likelihood was used).

implied: List. Model implied statistics.

vcov: List. Information about the variance matrix (vcov) of the model parameters.

test: List. Different test statistics.

h1: List. Information about the unrestricted h1 model (if available).

baseline: List. Information about a baseline model (often the independence model) (if available).

internal: List. For internal use only.

external: List. Empty slot to be used by add-on packages.

Methods

coef signature(object = "lavaan", type = "free"): Returns the estimates of the parameters in the model as a named numeric vector. If type="free", only the free parameters are returned. If type="user", all parameters listed in the parameter table are returned, including constrained and fixed parameters.

fitted.values signature(object = "lavaan"): Returns the implied moments of the model as a list with two elements (per group): cov for the implied covariance matrix, and mean for the implied mean vector. If only the covariance matrix was analyzed, the implied mean vector will be zero.

fitted signature(object = "lavaan"): an alias for fitted.values.

residuals signature(object = "lavaan", type="raw"): If type = "raw", this function returns the raw (= unscaled) difference between the observed and the expected (model-implied) summary statistics. If type = "cor", or type = "cor.bollen", the observed and model implied covariance matrices are first transformed to a correlation matrix (using cov2cor()), before the residuals are computed. If type = "cor.bentler", both the observed and model implied covariance matrices are rescaled by dividing the elements by the square roots of the corresponding variances of the observed covariance matrix. If type="normalized", the residuals are divided by the square root of the asymptotic variance of the corresponding summary statistic (the variance estimate depends on the choice for the se argument). Unfortunately, the corresponding normalized residuals are not entirely correct, and this option is only available for historical interest. If type="standardized", the residuals are divided by the square root of the asymptotic variance of these residuals. The resulting standardized residuals elements can be interpreted as z-scores. If type="standardized.mplus", the residuals are divided

by the square root of the asymptotic variance of these residuals. However, a simplified formula is used (see the Mplus reference below) which often results in negative estimates for the variances, resulting in many NA values for the standardized residuals.

- resid** signature(object = "lavaan"): an alias for residuals
- vcov** signature(object = "lavaan"): returns the covariance matrix of the estimated parameters.
- predict** signature(object = "lavaan"): compute factor scores for all cases that are provided in the data frame. For complete data only.
- anova** signature(object = "lavaan"): returns model comparison statistics. This method is just a wrapper around the function [lavTestLRT](#). If only a single argument (a fitted model) is provided, this model is compared to the unrestricted model. If two or more arguments (fitted models) are provided, the models are compared in a sequential order. Test statistics are based on the likelihood ratio test. For more details and further options, see the [lavTestLRT](#) page.
- update** signature(object = "lavaan", model, add, ..., evaluate=TRUE): update a fitted lavaan object and evaluate it (unless evaluate=FALSE). Note that we use the environment that is stored within the lavaan object, which is not necessarily the parent frame. The add argument is analogous to the one described in the [lavTestScore](#) page, and can be used to add parameters to the specified model rather than passing an entirely new model argument.
- nobs** signature(object = "lavaan"): returns the effective number of observations used when fitting the model. In a multiple group analysis, this is the sum of all observations per group.
- logLik** signature(object = "lavaan"): returns the log-likelihood of the fitted model, if maximum likelihood estimation was used. The [AIC](#) and [BIC](#) methods automatically work via [logLik\(\)](#).
- show** signature(object = "lavaan"): Print a short summary of the model fit
- summary** signature(object = "lavaan", header = TRUE, fit.measures = FALSE, estimates = TRUE, ci = FALSE, fmi = FALSE, standardized = FALSE, std.nox = FALSE, remove.step1 = TRUE, remove.unused = TRUE, cov.std = TRUE, rsquare = FALSE, modindices = FALSE, ci = FALSE, nd = 3L): Print a nice summary of the model estimates. If header = TRUE, the header section (including fit measures) is printed. If fit.measures = TRUE, additional fit measures are added to the header section. The related fm.args list allows to set options related to the fit measures. See [fitMeasures](#) for more details. If estimates = TRUE, print the parameter estimates section. If ci = TRUE, add confidence intervals to the parameter estimates section. If fmi = TRUE, add the fmi (fraction of missing information) column, if it is available. If standardized=TRUE or a character vector, the standardized solution is also printed (see [parameterEstimates](#)). Note that SEs and tests are still based on unstandardized estimates. Use [standardizedSolution](#) to obtain SEs and test statistics for standardized estimates. The std.nox argument is deprecated; the standardized argument allows "std.nox" solution to be specifically requested. If remove.step1, the parameters of the measurement part are not shown (only used when using sam().) If remove.unused, automatically added parameters that are fixed to their default (0 or 1) values are removed. If rsquare=TRUE, the R-Square values for the dependent variables in the model are printed. If efa = TRUE, EFA related information is printed. The related efa.args list allows to set options related to the EFA output. See [summary.efaList](#) for more details. If modindices=TRUE, modification indices are printed for all fixed parameters. The argument nd determines the number of digits after the decimal point to be printed (currently only in the parameter estimates section.) Historically, nothing was returned, but since 0.6-12, a list is returned of class lavaan.summary for which is print function is available.

References

Yves Rosseel (2012). lavaan: An R Package for Structural Equation Modeling. *Journal of Statistical Software*, 48(2), 1-36. doi:10.18637/jss.v048.i02

Standardized Residuals in Mplus. Document retrieved from URL <https://www.statmodel.com/download/StandardizedResiduals>

See Also

[cfa](#), [sem](#), [fitMeasures](#), [standardizedSolution](#), [parameterEstimates](#), [lavInspect](#), [modindices](#)

Examples

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

```
fit <- cfa(HS.model, data = HolzingerSwineford1939)
```

```
summary(fit, standardized = TRUE, fit.measures = TRUE, rsquare = TRUE)
fitted(fit)
coef(fit)
resid(fit, type = "normalized")
```

lavaanList

Fit List of Latent Variable Models

Description

Fit the same latent variable model, for a (potentially large) number of datasets.

Usage

```
lavaanList(model = NULL, dataList = NULL, dataFunction = NULL,
            dataFunction.args = list(), ndat = length(dataList), cmd = "lavaan",
            ..., store.slots = c("partable"), FUN = NULL, show.progress = FALSE,
            store.failed = FALSE, parallel = c("no", "multicore", "snow"),
            ncpus = max(1L, parallel::detectCores() - 1L), cl = NULL, iseed = NULL)
```

```
semList(model = NULL, dataList = NULL, dataFunction = NULL,
         dataFunction.args = list(), ndat = length(dataList),
         ..., store.slots = c("partable"), FUN = NULL, show.progress = FALSE,
         store.failed = FALSE, parallel = c("no", "multicore", "snow"),
         ncpus = max(1L, parallel::detectCores() - 1L), cl = NULL, iseed = NULL)
```

```
cfaList(model = NULL, dataList = NULL, dataFunction = NULL,
         dataFunction.args = list(), ndat = length(dataList),
         ..., store.slots = c("partable"), FUN = NULL, show.progress = FALSE,
         store.failed = FALSE, parallel = c("no", "multicore", "snow"),
         ncpus = max(1L, parallel::detectCores() - 1L), cl = NULL, iseed = NULL)
```

Arguments

<code>model</code>	A description of the user-specified model. Typically, the model is described using the lavaan model syntax. See model.syntax for more information. Alternatively, a parameter table (eg. the output of the <code>lavaanify()</code> function) is also accepted.
<code>dataList</code>	List. Each element contains a full data frame containing the observed variables used in the model.
<code>dataFunction</code>	Function. A function that generated a full data frame containing the observed variables used in the model. It can also be a matrix, if the columns are named.
<code>dataFunction.args</code>	List. Optional list of arguments that are passed to the <code>dataFunction</code> function.
<code>ndat</code>	Integer. The number of datasets that should be generated using the <code>dataFunction</code> function.
<code>cmd</code>	Character. Which command is used to run the sem models. The possible choices are "sem", "cfa" or "lavaan", determining how we deal with default options.
<code>...</code>	Other named arguments for lavaan function.
<code>store.slots</code>	Character vector. Which slots (from a lavaan object) should be stored for each dataset? The possible choices are "timing", "partable", "data", "samplestats", "vcov", "test", "optim", "h1", "loglik", or "implied". Finally, "all" selects all slots.
<code>FUN</code>	Function. A function which when applied to the lavaan object returns the information of interest.
<code>store.failed</code>	Logical. If TRUE, write (to <code>tempdir()</code>) the dataset and (if available) the fitted object when the estimation for a particular dataset somehow failed. This will allow posthoc inspection of the problem.
<code>parallel</code>	The type of parallel operation to be used (if any). If missing, the default is "no".
<code>ncpus</code>	Integer. The number of processes to be used in parallel operation: typically one would chose this to the number of available CPUs.
<code>cl</code>	An optional parallel or snow cluster for use if <code>parallel = "snow"</code> . If not supplied, a cluster on the local machine is created for the duration of the <code>lavaanList</code> call.
<code>iseed</code>	An integer to set the seed. Or NULL if no reproducible seeds are needed. To make this work, make sure the first <code>RNGkind()</code> element is "L'Ecuyer-CMRG". You can check this by typing <code>RNGkind()</code> in the console. You can set it by typing <code>RNGkind("L'Ecuyer-CMRG")</code> , before the <code>lavaanList</code> functions are called.
<code>show.progress</code>	If TRUE, show information for each dataset.

Value

An object of class [lavaanList](#), for which several methods are available, including a summary method.

See Also

class [lavaanList](#)

Examples

```
# The Holzinger and Swineford (1939) example
HS.model <- ' visual  =~ x1 + x2 + x3
            textual =~ x4 + x5 + x6
            speed   =~ x7 + x8 + x9 '

# a data generating function
generateData <- function() simulateData(HS.model, sample.nobs = 100)

set.seed(1234)
fit <- semList(HS.model, dataFunction = generateData, ndat = 5,
              store.slots = "partable")

# show parameter estimates, per dataset
coef(fit)
```

lavaanList-class

Class For Representing A List of (Fitted) Latent Variable Models

Description

The `lavaanList` class represents a collection of (fitted) latent variable models, for a (potentially large) number of datasets. It contains information about the model (which is always the same), and for every dataset a set of (user-specified) slots from a regular lavaan object.

Objects from the Class

Objects can be created via the `cfaList`, `semList`, or `lavaanList` functions.

Slots

version: The lavaan package version used to create this objects

call: The function call as returned by `match.call()`.

Options: Named list of options that were provided by the user, or filled-in automatically.

ParTable: Named list describing the model parameters. Can be coerced to a `data.frame`. In the documentation, this is called the ‘parameter table’.

pta: Named list containing parameter table attributes.

Data: Object of internal class "Data": information about the data.

Model: Object of internal class "Model": the internal (matrix) representation of the model

meta: List containing additional flags. For internal use only.

timingList: List. Timing slot per dataset.

ParTableList: List. ParTable slot per dataset.

DataList: List. Data slot per dataset.

SampleStatsList: List. SampleStats slot per dataset.

CacheList: List. Cache slot per dataset.
vcovList: List. vcov slot per dataset.
testList: List. test slot per dataset.
optimList: List. optim slot per dataset.
impliedList: List. implied slot per dataset.
h1List: List. h1 slot per dataset.
loglikList: List. loglik slot per dataset.
baselineList: List. baseline slot per dataset.
funList: List. fun slot per dataset.
internalList: List. internal slot per dataset.
external: List. Empty slot to be used by add-on packages.

Methods

coef signature(object = "lavaanList", type = "free"): Returns the estimates of the parameters in the model as the columns in a matrix; each column corresponds to a different dataset. If type="free", only the free parameters are returned. If type="user", all parameters listed in the parameter table are returned, including constrained and fixed parameters.

summary signature(object = "lavaanList", header = TRUE, estimates = TRUE, nd = 3L): Print a summary of the collection of fitted models. If header = TRUE, the header section is printed. If estimates = TRUE, print the parameter estimates section. The argument nd determines the number of digits after the decimal point to be printed (currently only in the parameter estimates section.) Nothing is returned (use parameterEstimates or another extractor function to extract information from this object).

See Also

[cfaList](#), [semList](#), [lavaanList](#)

lavCor

Polychoric, polyserial and Pearson correlations

Description

Fit an unrestricted model to compute polychoric, polyserial and/or Pearson correlations.

Usage

```

lavCor(object, ordered = NULL, group = NULL, missing = "listwise",
       ov.names.x = NULL, sampling.weights = NULL,
       se = "none", test = "none",
       estimator = "two.step", baseline = FALSE, ...,
       cor.smooth = FALSE, cor.smooth.tol = 1e-04, output = "cor")
  
```

Arguments

object	Either a <code>data.frame</code> , or an object of class <code>lavaan</code> . If the input is a <code>data.frame</code> , and some variables are declared as ordered factors, <code>lavaan</code> will treat them as ordinal variables.
ordered	Character vector. Only used if object is a <code>data.frame</code> . Treat these variables as ordered (ordinal) variables. Importantly, all other variables will be treated as numeric (unless they are declared as ordered in the original data frame.)
group	Only used if object is a <code>data.frame</code> . Specify a grouping variable.
missing	If "listwise", cases with missing values are removed listwise from the data frame. If "direct" or "ml" or "fiml" and the estimator is maximum likelihood, an EM algorithm is used to estimate the unrestricted covariance matrix (and mean vector). If "pairwise", pairwise deletion is used. If "default", the value is set depending on the estimator and the mimic option.
sampling.weights	Only used if object is a <code>data.frame</code> . Specify a variable containing sampling weights.
ov.names.x	Only used if object is a <code>data.frame</code> . Specify variables that need to be treated as exogenous. Only used if at least one variable is declared as ordered.
se	Only used if output (see below) contains standard errors. See lavOptions for possible options.
test	Only used if output is "fit" or "lavaan". See lavOptions for possible options.
estimator	If "none" or "two.step" or "two.stage", only starting values are computed for the correlations (and thresholds), without any further estimation. If all variables are continuous, the starting values are the sample covariances (converted to correlations if output = "cor"). If at least one variable is ordered, the thresholds are computed using univariate information only. The polychoric and/or polyserial correlations are computed in a second stage, keeping the values of the thresholds constant. If an estimator (other than "two.step" or "two.stage") is specified (for example estimator = "PML"), these starting values are further updated by fitting the unrestricted model using the chosen estimator. See the lavaan function for alternative estimators.
baseline	Only used if output is "fit" or "lavaan". If TRUE, a baseline model is also estimated. Note that the test argument should also be set to a value other than "none".
...	Optional parameters that are passed to the lavaan function.
cor.smooth	Logical. Only used if output = "cor". If TRUE, ensure the resulting correlation matrix is positive definite. The following simple method is used: an eigenvalue decomposition is computed; then, eigenvalues smaller than <code>cor.smooth.tol</code> are set to be equal to <code>cor.smooth.tol</code> , before the matrix is again reconstructed. Finally, the matrix (which may no longer have unit diagonal elements) is converted to a correlation matrix using <code>cov2cor</code> .
cor.smooth.tol	Numeric. Smallest eigenvalue used when reconstructing the correlation matrix after an eigenvalue decomposition.

output If "cor", the function returns the correlation matrix only. If "cov", the function returns the covariance matrix (this only makes a difference if at least one variable is numeric). If "th" or "thresholds", only the thresholds are returned. If "sampstat", the output equals the result of `lavInspect(fit, "sampstat")` where `fit` is the unrestricted model. If "est" or "pe" or "parameterEstimates", the output equals the result of `parameterEstimates(fit)`. Finally, if output is "fit" or "lavaan", the function returns an object of class `lavaan`.

Details

This function is a wrapper around the `lavaan` function, but where the model is defined as the unrestricted model. The following free parameters are included: all covariances/correlations among the variables, variances for continuous variables, means for continuous variables, thresholds for ordered variables, and if exogenous variables are included (`ov.names.x` is not empty) while some variables are ordered, also the regression slopes enter the model.

Value

By default, if `output = "cor"` or `output = "cov"`, a symmetric matrix (of class `"lavaan.matrix.symmetric"`, which only affects the way the matrix is printed). If `output = "th"`, a named vector of thresholds. If `output = "fit"` or `output = "lavaan"`, an object of class `lavaan`.

References

Olsson, U. (1979). Maximum likelihood estimation of the polychoric correlation coefficient. *Psychometrika*, 44(4), 443-460.

Olsson, U., Drasgow, F., & Dorans, N. J. (1982). The polyserial correlation coefficient. *Psychometrika*, 47(3), 337-347.

See Also

[lavaan](#)

Examples

```
# Holzinger and Swineford (1939) example
HS9 <- HolzingerSwineford1939[,c("x1", "x2", "x3", "x4", "x5",
                                "x6", "x7", "x8", "x9")]

# Pearson correlations
lavCor(HS9)

# ordinal version, with three categories
HS9ord <- as.data.frame( lapply(HS9, cut, 3, labels = FALSE) )

# polychoric correlations, two-stage estimation
lavCor(HS9ord, ordered=names(HS9ord))

# thresholds only
lavCor(HS9ord, ordered=names(HS9ord), output = "th")
```

```
# polychoric correlations, with standard errors
lavCor(HS9ord, ordered=names(HS9ord), se = "standard", output = "est")

# polychoric correlations, full output
fit.un <- lavCor(HS9ord, ordered=names(HS9ord), se = "standard",
                 output = "fit")
summary(fit.un)
```

lavExport

lavaan Export

Description

Export a fitted lavaan object to an external program.

Usage

```
lavExport(object, target = "lavaan", prefix = "sem", dir.name = "lavExport",
          export = TRUE)
```

Arguments

object	An object of class lavaan .
target	The target program. Current options are "lavaan" and "Mplus".
prefix	The prefix used to create the input files; the name of the input file has the pattern 'prefix dot target dot in'; the name of the data file has the pattern 'prefix dot target dot raw'.
dir.name	The directory name (including a full path) where the input files will be written.
export	If TRUE, the files are written to the output directory (dir.name). If FALSE, only the syntax is generated as a character string.

Details

This function was mainly created to quickly generate an Mplus syntax file to compare the results between Mplus and lavaan. The target "lavaan" can be useful to create a full model syntax as needed for the lavaan() function. More targets (perhaps for LISREL or EQS) will be added in future releases.

Value

If export = TRUE, a directory (called lavExport by default) will be created, typically containing a data file, and an input file so that the same analysis can be run using an external program. If export = FALSE, a character string containing the model syntax only for the target program.

See Also

[lavaanify](#), [mplus2lavaan](#)

Examples

```

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

fit <- cfa(HS.model, data=HolzingerSwineford1939)
out <- lavExport(fit, target = "Mplus", export=FALSE)
cat(out)

```

lavInspect

*Inspect or extract information from a fitted lavaan object***Description**

The `lavInspect()` and `lavTech()` functions can be used to inspect/extract information that is stored inside (or can be computed from) a fitted lavaan object. Note: the (older) `inspect()` function is now simply a shortcut for `lavInspect()` with default arguments.

Usage

```

lavInspect(object, what = "free", add.labels = TRUE, add.class = TRUE,
           list.by.group = TRUE,
           drop.list.single.group = TRUE)

```

```

lavTech(object, what = "free", add.labels = FALSE, add.class = FALSE,
        list.by.group = FALSE,
        drop.list.single.group = FALSE)

```

```

inspect(object, what = "free", ...)

```

Arguments

<code>object</code>	An object of class <code>lavaan</code> .
<code>what</code>	Character. What needs to be inspected/extracted? See Details for a full list. Note: the <code>what</code> argument is not case-sensitive (everything is converted to lower case.)
<code>add.labels</code>	If TRUE, variable names are added to the vectors and/or matrices.
<code>add.class</code>	If TRUE, vectors are given the <code>'lavaan.vector'</code> class; matrices are given the <code>'lavaan.matrix'</code> class, and symmetric matrices are given the <code>'lavaan.matrix.symmetric'</code> class. This only affects the way they are printed on the screen.
<code>list.by.group</code>	Logical. Only used when the output are model matrices. If TRUE, the model matrices are nested within groups. If FALSE, a flattened list is returned containing all model matrices, with repeated names for multiple groups.
<code>drop.list.single.group</code>	If FALSE, the results are returned as a list, where each element corresponds to a group (even if there is only a single group). If TRUE, the list will be unlisted if there is only a single group.
<code>...</code>	Additional arguments. Not used by <code>lavaan</code> , but by other packages.

Details

The `lavInspect()` and `lavTech()` functions only differ in the way they return the results. The `lavInspect()` function will prettify the output by default, while the `lavTech()` will not attempt to prettify the output by default. The (older) `inspect()` function is a simplified version of `lavInspect()` with only the first two arguments.

Below is a list of possible values for the `what` argument, organized in several sections:

Model matrices:

`"free"`: A list of model matrices. The non-zero integers represent the free parameters. The numbers themselves correspond to the position of the free parameter in the parameter vector. This determines the order of the model parameters in the output of for example `coef()` and `vcov()`.

`"partable"`: A list of model matrices. The non-zero integers represent both the fixed parameters (for example, factor loadings fixed at 1.0), and the free parameters if we ignore any equality constraints. They correspond with all entries (fixed or free) in the parameter table. See [parTable](#).

`"se"`: A list of model matrices. The non-zero numbers represent the standard errors for the free parameters in the model. If two parameters are constrained to be equal, they will have the same standard error for both parameters. Aliases: `"std.err"` and `"standard.errors"`.

`"start"`: A list of model matrices. The values represent the starting values for all model parameters. Alias: `"starting.values"`.

`"est"`: A list of model matrices. The values represent the estimated model parameters. Aliases: `"estimates"`, and `"x"`.

`"dx.free"`: A list of model matrices. The values represent the gradient (first derivative) values of the model parameters. If two parameters are constrained to be equal, they will have the same gradient value.

`"dx.all"`: A list of model matrices. The values represent the first derivative with respect to all possible matrix elements. Currently, this is only available when the estimator is "ML" or "GLS".

`"std"`: A list of model matrices. The values represent the (completely) standardized model parameters (the variances of both the observed and the latent variables are set to unity). Aliases: `"std.all"`, `"standardized"`.

`"std.lv"`: A list of model matrices. The values represent the standardized model parameters (only the variances of the latent variables are set to unity.)

`"std.nox"`: A list of model matrices. The values represent the (completely) standardized model parameters (the variances of both the observed and the latent variables are set to unity; however, the variances of any observed exogenous variables are not set to unity; hence no-x.)

Information about the data:

`"data"`: A matrix containing the observed variables that have been used to fit the model. No column/row names are provided. Column names correspond to the output of `lavNames(object)`, while the rows correspond to the output of `lavInspect(object, "case.idx")`.

`"ordered"`: A character vector. The ordered variables.

`"nobs"`: Integer vector. The number of observations in each group that were used in the analysis.

- "norig": Integer vector. The original number of observations in each group.
- "ntotal": Integer. The total number of observations that were used in the analysis. If there is just a single group, this is the same as the "nobs" option; if there are multiple groups, this is the sum of the "nobs" numbers for each group.
- "case.idx": Integer vector. The case/observation numbers that were used in the analysis. In the case of multiple groups: a list of numbers.
- "empty.idx": The case/observation numbers of those cases/observations that contained missing values only (at least for the observed variables that were included in the model). In the case of multiple groups: a list of numbers.
- "patterns": A binary matrix. The rows of the matrix are the missing data patterns where 1 and 0 denote non-missing and missing values for the corresponding observed variables respectively (or TRUE and FALSE if lavTech() is used.) If the data is complete (no missing values), there will be only a single pattern. In the case of multiple groups: a list of pattern matrices.
- "coverage": A symmetric matrix where each element contains the proportion of observed data-points for the corresponding pair of observed variables. In the case of multiple groups: a list of coverage matrices.
- "group": A character string. The group variable in the data.frame (if any).
- "ngroups": Integer. The number of groups.
- "group.label": A character vector. The group labels.
- "level.label": A character vector. The level labels.
- "cluster": A character vector. The cluster variable(s) in the data.frame (if any).
- "nlevels": Integer. The number of levels.
- "nclusters": Integer. The number of clusters that were used in the analysis.
- "ncluster.size": Integer. The number of different cluster sizes.
- "cluster.size": Integer vector. The number of observations within each cluster. For multigroup multilevel models, a list of integer vectors, indicating cluster sizes within each group.
- "cluster.id": Integer vector. The cluster IDs identifying the clusters. For multigroup multilevel models, a list of integer vectors, indicating cluster IDs within each group.
- "cluster.idx": Integer vector. The cluster index for each observation. The cluster index ranges from 1 to the number of clusters. For multigroup multilevel models, a list of integer vectors, indicating cluster indices within each group.
- "cluster.label": Integer vector. The cluster ID for each observation. For multigroup multilevel models, a list of integer vectors, indicating the cluster ID for each observation within each group.
- "cluster.sizes": Integer vector. The different cluster sizes that were used in the analysis. For multigroup multilevel models, a list of integer vectors, indicating the different cluster sizes within each group.
- "average.cluster.size": Integer. The average cluster size (using the formula $s = (N^2 - \text{sum}(\text{cluster.size}^2)) / (N * (\text{nclusters} - 1L))$). For multigroup multilevel models, a list containing the average cluster size per group.

Observed sample statistics:

"smpstat": Observed sample statistics. Aliases: "obs", "observed", "smp", "sample", "samplestatistics".

Since 0.6-3, we always check if an h1 slot is available (the estimates for the unrestricted model); if present, we extract the sample statistics from this slot. This implies that if variables are continuous, and missing = "ml" (or "fiml"), we return the covariance matrix (and mean vector) as computed by the EM algorithm under the unrestricted (h1) model. If the h1 is not present (perhaps, because the model was fitted with h1 = FALSE), we return the sample statistics from the SampleStats slot. Here, pairwise deletion is used for the elements of the covariance matrix (or correlation matrix), and listwise deletion for all univariate statistics (means, intercepts and thresholds).

"smpstat.h1": Deprecated. Do not use any longer.

"wls.obs": The observed sample statistics (covariance elements, intercepts/thresholds, etc.) in a single vector.

"wls.v": The weight vector as used in weighted least squares estimation.

"gamma": N times the asymptotic variance matrix of the sample statistics. Alias: "smpstat.nacov".

Model features:

"meanstructure": Logical. TRUE if a meanstructure was included in the model.

"categorical": Logical. TRUE if categorical endogenous variables were part of the model.

"fixed.x": Logical. TRUE if the exogenous x-covariates are treated as fixed.

"parameterization": Character. Either "delta" or "theta".

Model-implied sample statistics:

"implied": The model-implied summary statistics. Alias: "fitted", "expected", "exp".

"resid": The difference between observed and model-implied summary statistics. Alias: "residuals", "residual", "res".

"cov.lv": The model-implied variance-covariance matrix of the latent variables. Alias: "veta" [for V(eta)].

"cor.lv": The model-implied correlation matrix of the latent variables.

"mean.lv": The model-implied mean vector of the latent variables. Alias: "eeta" [for E(eta)].

"cov.ov": The model-implied variance-covariance matrix of the observed variables. Aliases: "sigma", "sigma.hat".

"cor.ov": The model-implied correlation matrix of the observed variables.

"mean.ov": The model-implied mean vector of the observed variables. Aliases: "mu", "mu.hat".

"cov.all": The model-implied variance-covariance matrix of both the observed and latent variables.

"cor.all": The model-implied correlation matrix of both the observed and latent variables.

"th": The model-implied thresholds. Alias: "thresholds".

"wls.est": The model-implied sample statistics (covariance elements, intercepts/thresholds, etc.) in a single vector.

"vy": The model-implied unconditional variances of the observed variables.

"rsquare": The R-square value for all endogenous variables. Aliases: "r-square", "r2".

"fs.determinacy": The factor determinacies (based on regression factor scores). They represent the (estimated) correlation between the factor scores and the latent variables scores.

"fs.reliability": The factor reliabilities (based on regression factor scores). They are the square of the factor determinacies.

"fs.determinacy.Bartlett": The factor determinacies (based on Bartlett factor scores). They represent the (estimated) correlation between the factor scores and the latent variables scores.

"fs.reliability.Bartlett": The factor reliabilities (based on Bartlett factor scores). They are the square of the factor determinacies.

Diagnostics:

"mdist2.fs": The squared Mahalanobis distances for the (Bartlett) factor scores.

"mdist.fs": The Mahalanobis distances for the (Bartlett) factor scores.

"mdist2.resid": The squared Mahalanobis distances for the (Bartlett-based) casewise residuals.

"mdist.resid": The Mahalanobis distances for the (Bartlett-based) casewise residuals.

Optimizer information:

"converged": Logical. TRUE if the optimizer has converged; FALSE otherwise.

"iterations": Integer. The number of iterations used by the optimizer.

"optim": List. All available information regarding the optimization results.

"npar": Integer. Number of free parameters (ignoring constraints).

Gradient, Hessian, observed, expected and first.order information matrices:

"gradient": Numeric vector containing the first derivatives of the discrepancy function with respect to the (free) model parameters.

"hessian": Matrix containing the second derivatives of the discrepancy function with respect to the (free) model parameters.

"information": Matrix containing either the observed or the expected information matrix (depending on the information option of the fitted model). This is unit-information, not total-information.

"information.expected": Matrix containing the expected information matrix for the free model parameters.

"information.observed": Matrix containing the observed information matrix for the free model parameters.

"information.first.order": Matrix containing the first.order information matrix for the free model parameters. This is the outer product of the gradient elements (the first derivative of the discrepancy function with respect to the (free) model parameters). Alias: "first.order".

"augmented.information": Matrix containing either the observed or the expected augmented (or bordered) information matrix (depending on the information option of the fitted model). Only relevant if constraints have been used in the model.

"augmented.information.expected": Matrix containing the expected augmented (or bordered) information matrix. Only relevant if constraints have been used in the model.

- "augmented.information.observed": Matrix containing the observed augmented (or bordered) information matrix. Only relevant if constraints have been used in the model.
- "augmented.information.first.order": Matrix containing the first.order augmented (or bordered) information matrix. Only relevant if constraints have been used in the model.
- "inverted.information": Matrix containing either the observed or the expected inverted information matrix (depending on the information option of the fitted model).
- "inverted.information.expected": Matrix containing the inverted expected information matrix for the free model parameters.
- "inverted.information.observed": Matrix containing the inverted observed information matrix for the free model parameters.
- "inverted.information.first.order": Matrix containing the inverted first.order information matrix for the free model parameters.
- "h1.information": Matrix containing either the observed, expected or first.order information matrix (depending on the information option of the fitted model) of the unrestricted h1 model. This is unit-information, not total-information.
- "h1.information.expected": Matrix containing the expected information matrix for the unrestricted h1 model.
- "h1.information.observed": Matrix containing the observed information matrix for the unrestricted h1 model.
- "h1.information.first.order": Matrix containing the first.order information matrix for the the unrestricted h1 model. Alias: "h1.first.order".

Variance covariance matrix of the model parameters:

- "vcov": Matrix containing the variance covariance matrix of the estimated model parameters.
- "vcov.std.all": Matrix containing the variance covariance matrix of the standardized estimated model parameters. Standardization is done with respect to both observed and latent variables.
- "vcov.std.lv": Matrix containing the variance covariance matrix of the standardized estimated model parameters. Standardization is done with respect to the latent variables only.
- "vcov.std.nox": Matrix containing the variance covariance matrix of the standardized estimated model parameters. Standardization is done with respect to both observed and latent variables, but ignoring any exogenous observed covariates.
- "vcov.def": Matrix containing the variance covariance matrix of the user-defined (using the := operator) parameters.
- "vcov.def.std.all": Matrix containing the variance covariance matrix of the standardized user-defined parameters. Standardization is done with respect to both observed and latent variables.
- "vcov.def.std.lv": Matrix containing the variance covariance matrix of the standardized user-defined parameters. Standardization is done with respect to the latent variables only.
- "vcov.def.std.nox": Matrix containing the variance covariance matrix of the standardized user-defined parameters. Standardization is done with respect to both observed and latent variables, but ignoring any exogenous observed covariates.
- "vcov.def.joint": Matrix containing the joint variance covariance matrix of both the estimated model parameters and the defined (using the := operator) parameters.

"vcov.def.joint.std.all": Matrix containing the joint variance covariance matrix of both the standardized model parameters and the user-defined parameters. Standardization is done with respect to both observed and latent variables.

"vcov.def.joint.std.lv": Matrix containing the joint variance covariance matrix of both the standardized model parameters and the user-defined parameters. Standardization is done with respect to the latent variables only.

"vcov.def.joint.std.nox": Matrix containing the joint variance covariance matrix of both the standardized model parameters and the user-defined parameters. Standardization is done with respect to both observed and latent variables, but ignoring any exogenous observed covariates.

Miscellaneous:

"coef.boot": Matrix containing estimated model parameters for for each bootstrap sample. Only relevant when bootstrapping was used.

"UGamma": Matrix containing the product of 'U' and 'Gamma' matrices as used by the Satorra-Bentler correction. The trace of this matrix, divided by the degrees of freedom, gives the scaling factor.

"UfromUGamma": Matrix containing the 'U' matrix as used by the Satorra-Bentler correction. Alias: "U".

"list": The parameter table. The same output as given by parTable().

"fit": The fit measures. Aliases: "fitmeasures", "fit.measures", "fit.indices". The same output as given by fitMeasures().

"mi": The modification indices. Alias: "modindices", "modification.indices". The same output as given by modindices().

"loglik.casewise": Vector containing the casewise loglikelihood contributions. Only available if estimator = "ML".

"options": List. The option list.

"call": List. The call as returned by match.call, coerced to a list.

"timing": List. The timing (in milliseconds) of various lavaan subprocedures.

"test": List. All available information regarding the (goodness-of-fit) test statistic(s).

"baseline.test": List. All available information regarding the (goodness-of-fit) test statistic(s) of the baseline model.

"baseline.parTable": Data.frame. The parameter table of the (internal) baseline model.

"post.check": Post-fitting check if the solution is admissible. A warning is raised if negative variances are found, or if either lavInspect(fit, "cov.lv") or lavInspect(fit, "theta") return a non-positive definite matrix.

"zero.cell.tables": List. List of bivariate frequency tables where at least one cell is empty.

"version": The lavaan version number that was used to construct the fitted lavaan object.

See Also

[lavaan](#)

Examples

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

fit <- cfa(HS.model, data = HolzingerSwineford1939, group = "school")

# extract information
lavInspect(fit, "sampstat")
lavTech(fit, "sampstat")
```

lavListInspect	<i>Inspect or extract information from a lavaanList object</i>
----------------	--

Description

The `lavListInspect()` and `lavListTech()` functions can be used to inspect/extract information that is stored inside (or can be computed from) a `lavaanList` object.

Usage

```
lavListInspect(object, what = "free", add.labels = TRUE,
               add.class = TRUE, list.by.group = TRUE,
               drop.list.single.group = TRUE)
```

```
lavListTech(object, what = "free", add.labels = FALSE,
             add.class = FALSE, list.by.group = FALSE,
             drop.list.single.group = FALSE)
```

Arguments

<code>object</code>	An object of class <code>lavaanList</code> .
<code>what</code>	Character. What needs to be inspected/extracted? See Details for a full list. Note: the <code>what</code> argument is not case-sensitive (everything is converted to lower case.)
<code>add.labels</code>	If TRUE, variable names are added to the vectors and/or matrices.
<code>add.class</code>	If TRUE, vectors are given the ‘ <code>lavaan.vector</code> ’ class; matrices are given the ‘ <code>lavaan.matrix</code> ’ class, and symmetric matrices are given the ‘ <code>lavaan.matrix.symmetric</code> ’ class. This only affects the way they are printed on the screen.
<code>list.by.group</code>	Logical. Only used when the output are model matrices. If TRUE, the model matrices are nested within groups. If FALSE, a flattened list is returned containing all model matrices, with repeated names for multiple groups.
<code>drop.list.single.group</code>	If FALSE, the results are returned as a list, where each element corresponds to a group (even if there is only a single group.) If TRUE, the list will be unlisted if there is only a single group.

Details

The `lavListInspect()` and `lavListTech()` functions only differ in the way they return the results. The `lavListInspect()` function will prettify the output by default, while the `lavListTech()` will not attempt to prettify the output by default.

Below is a list of possible values for the `what` argument, organized in several sections:

Model matrices:

`"free"`: A list of model matrices. The non-zero integers represent the free parameters. The numbers themselves correspond to the position of the free parameter in the parameter vector. This determines the order of the model parameters in the output of for example `coef()` and `vcov()`.

`"partable"`: A list of model matrices. The non-zero integers represent both the fixed parameters (for example, factor loadings fixed at 1.0), and the free parameters if we ignore any equality constraints. They correspond with all entries (fixed or free) in the parameter table. See [parTable](#).

`"start"`: A list of model matrices. The values represent the starting values for all model parameters. Alias: `"starting.values"`.

Information about the data (including missing patterns):

`"group"`: A character string. The group variable in the `data.frame` (if any).

`"ngroups"`: Integer. The number of groups.

`"group.label"`: A character vector. The group labels.

`"level.label"`: A character vector. The level labels.

`"cluster"`: A character vector. The cluster variable(s) in the `data.frame` (if any).

`"nlevels"`: Integer. The number of levels.

`"ordered"`: A character vector. The ordered variables.

`"nobs"`: Integer vector. The number of observations in each group that were used in the analysis (in each dataset).

`"norig"`: Integer vector. The original number of observations in each group (in each dataset).

`"ntotal"`: Integer. The total number of observations that were used in the analysis. If there is just a single group, this is the same as the `"nobs"` option; if there are multiple groups, this is the sum of the `"nobs"` numbers for each group (in each dataset).

Model features:

`"meanstructure"`: Logical. TRUE if a meanstructure was included in the model.

`"categorical"`: Logical. TRUE if categorical endogenous variables were part of the model.

`"fixed.x"`: Logical. TRUE if the exogenous x -covariates are treated as fixed.

`"parameterization"`: Character. Either `"delta"` or `"theta"`.

`"list"`: The parameter table. The same output as given by `parTable()`.

`"options"`: List. The option list.

`"call"`: List. The call as returned by `match.call`, coerced to a list.

See Also[lavaanList](#)**Examples**

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

# a data generating function
generateData <- function() simulateData(HS.model, sample.nobs = 100)

set.seed(1234)
fit <- semList(HS.model, dataFunction = generateData, ndat = 5,
              store.slots = "partable")

# extract information
lavListInspect(fit, "free")
lavListTech(fit, "free")
```

lavMatrixRepresentation

lavaan matrix representation

Description

Extend the parameter table with a matrix representation.

Usage

```
lavMatrixRepresentation(partable, representation = "LISREL",
                       add.attributes = FALSE, as.data.frame. = TRUE)
```

Arguments

partable	A lavaan parameter table (as extracted by the parTable function, or generated by the lavPartable function).
representation	Character. The matrix representation style. Currently, only the all-y version of the LISREL representation is supported.
add.attributes	Logical. If TRUE, additional information about the model matrix representation is added as attributes.
as.data.frame.	Logical. If TRUE, the extended parameter table is returned as a data.frame.

Value

A list or a data.frame containing the original parameter table, plus three columns: a "mat" column containing matrix names, and a "row" and "col" column for the row and column indices of the model parameters in the model matrices.

See Also

[lavParTable](#), [parTable](#)

Examples

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

fit <- cfa(HS.model, data=HolzingerSwineford1939)

# extract partable
partable <- parTable(fit)

# add matrix representation (and show only a few columns)
lavMatrixRepresentation(partable)[,c("lhs", "op", "rhs", "mat", "row", "col")]
```

lavNames

lavaan Names

Description

Extract variables names from a fitted lavaan object.

Usage

```
lavNames(object, type = "ov", ...)
```

Arguments

object	An object of class lavaan .
type	Character. The type of variables whose names should be extracted. See details for a complete list.
...	Additional selection variables. For example "group = 2L" (in a multiple-group analysis) only considers the variables included in the model for the second group.

Details

The order of the variable names, as returned by `lavNames` determines the order in which the variables are listed in the parameter table, and therefore also in the summary output.

The following variable types are available:

- "ov": observed variables
- "ov.x": (pure) exogenous observed variables (no mediators)
- "ov.nox": non-exogenous observed variables
- "ov.model": modelled observed variables (joint vs conditional)

- "ov.y": (pure) endogenous variables (dependent only) (no mediators)
- "ov.num": numeric observed variables
- "ov.ord": ordinal observed variables
- "ov.ind": observed indicators of latent variables
- "ov.orphan": lonely observed variables (only intercepts/variances appear in the model syntax)
- "ov.interaction": interaction terms (defined by the colon operator)
- "th": threshold names ordinal variables only
- "th.mean": threshold names ordinal + numeric variables (if any)
- "lv": latent variables
- "lv.regular": latent variables (defined by =~ only)
- "lv.formative": latent variables (defined by <~ only)
- "lv.x": (pure) exogenous variables
- "lv.y": (pure) endogenous variables
- "lv.nox": non-exogenous latent variables
- "lv.nonnormal": latent variables with non-normal indicators
- "lv.interaction": interaction terms at the latent level
- "eqs.y": variables that appear as dependent variables in a regression formula (but not indicators of latent variables)
- "eqs.x": variables that appear as independent variables in a regression formula

See Also

[lavaanify](#), [parTable](#)

Examples

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

fit <- cfa(HS.model, data=HolzingerSwineford1939)
lavNames(fit, "ov")
```

lavOptions

*lavaan Options***Description**

Show the default options used by the `lavaan()` function. The options can be changed by passing 'name = value' arguments to the `lavaan()` function call, where they will be added to the '...' argument.

Usage

```
lavOptions(x = NULL, default = NULL, mimic = "lavaan")
```

Arguments

<code>x</code>	Character. A character string holding an option name, or a character string vector holding multiple option names. All option names are converted to lower case.
<code>default</code>	If a single option is specified but not available, this value is returned.
<code>mimic</code>	Not used for now.

Details

This is the full list of options that are accepted by the `lavaan()` function, organized in several sections:

Model features (always available):

`meanstructure`: If TRUE, the means of the observed variables enter the model. If "default", the value is set based on the user-specified model, and/or the values of other arguments.

`int.ov.free`: If FALSE, the intercepts of the observed variables are fixed to zero.

`int.lv.free`: If FALSE, the intercepts of the latent variables are fixed to zero.

`conditional.x`: If TRUE, we set up the model conditional on the exogenous 'x' covariates; the model-implied sample statistics only include the non-x variables. If FALSE, the exogenous 'x' variables are modeled jointly with the other variables, and the model-implied statistics reflect both sets of variables. If "default", the value is set depending on the estimator, and whether or not the model involves categorical endogenous variables.

`fixed.x`: If TRUE, the exogenous 'x' covariates are considered fixed variables and the means, variances and covariances of these variables are fixed to their sample values. If FALSE, they are considered random, and the means, variances and covariances are free parameters. If "default", the value is set depending on the mimic option.

`orthogonal`: If TRUE, all covariances among latent variables are set to zero.

`orthogonal.y`: If TRUE, all covariances among endogenous latent variables only are set to zero.

`orthogonal.x`: If TRUE, all covariances among exogenous latent variables only are set to zero.

- `std.lv`: If TRUE, the metric of each latent variable is determined by fixing their (residual) variances to 1.0. If FALSE, the metric of each latent variable is determined by fixing the factor loading of the first indicator to 1.0. If there are multiple groups, `std.lv = TRUE` and "loadings" is included in the `group.equal` argument, then only the latent variances of the first group will be fixed to 1.0, while the latent variances of other groups are set free.
- `effect.coding`: Can be logical or character string. If logical and TRUE, this implies `effect.coding = c("loadings", "intercepts")`. If logical and FALSE, it is set equal to the empty string. If "loadings" is included, equality constraints are used so that the average of the factor loadings (per latent variable) equals 1. Note that this should not be used together with `std.lv = TRUE`. If "intercepts" is included, equality constraints are used so that the sum of the intercepts (belonging to the indicators of a single latent variable) equals zero. As a result, the latent mean will be freely estimated and usually equal the average of the means of the involved indicators.
- `ceq.simple`: Logical. If TRUE, and no other general (equality or inequality) constraints are used in the model, simple equality constraints are represented in the parameter table as duplicated free parameters (instead of extra rows with `op = "=="`).
- `parameterization`: Currently only used if data is categorical. If "delta", the delta parameterization is used. If "theta", the theta parameterization is used.
- `correlation`: Only used for (single-level) continuous data. If TRUE, analyze a correlation matrix (instead of a (co)variance matrix). This implies that the residual observed variances are no longer free parameters. Instead, they are set to values to ensure the model-implied variances are unity. This also affects the standard errors. The only available estimators are GLS and WLS, which produce correct standard errors and a correct test statistic under normal and non-normal conditions respectively. Always assuming `fixed.x = FALSE` and `conditional.x = FALSE` (for now).

Model features (only available for the `lavaan()` function):

- `auto.fix.first`: If TRUE, the factor loading of the first indicator is set to 1.0 for every latent variable.
- `auto.fix.single`: If TRUE, the residual variance (if included) of an observed indicator is set to zero if it is the only indicator of a latent variable.
- auto.var** If TRUE, the (residual) variances of both observed and latent variables are set free.
- `auto.cov.lv.x`: If TRUE, the covariances of exogenous latent variables are included in the model and set free.
- `auto.cov.y`: If TRUE, the covariances of dependent variables (both observed and latent) are included in the model and set free.
- `auto.th`: If TRUE, thresholds for limited dependent variables are included in the model and set free.
- `auto.delta`: If TRUE, response scaling parameters for limited dependent variables are included in the model and set free.
- `auto.efa`: If TRUE, the necessary constraints are imposed to make the (unrotated) exploratory factor analysis blocks identifiable: for each block, factor variances are set to 1, factor covariances are constrained to be zero, and factor loadings are constrained to follow an echelon pattern.

Data options:

- std.ov:** If TRUE, observed variables are standardized before entering the analysis. By default, these are only the non-exogenous observed variables, unless `fixed.x = FALSE`. Use this option with caution; it can be used to test if (for example) nonconvergence was due to scaling issues. But this is still a covariance based analysis, in the sense that no constraints are involved (to ensure the model-implied (co)variance matrix has unit variances), and the standard errors still assume that the input was unstandardized. See also the `correlation` option.
- missing:** The default setting is "listwise": all cases with missing values are removed listwise from the data before the analysis starts. This is only valid if the data are missing completely at random (MCAR). Therefore, it may not be the optimal choice, but it can be useful for a first run. If the estimator belongs to the ML family, another option is "ml" (alias: "fiml" or "direct"). This corresponds to the so-called full information maximum likelihood approach (fiml), where we compute the likelihood case by case, using all available data from that case. Note that if the model contains exogenous observed covariates, and `fixed.x = TRUE` (the default), all cases with any missing values on these covariates will be deleted first. The option "ml.x" (alias: "fiml.x" or "direct.x") is similar to "ml", but does not delete any cases with missing values for the exogenous covariates, even if `fixed.x = TRUE`. (Note: all lavaan versions < 0.6 used "ml.x" instead of "ml"). If you wish to use multiple imputation, you need to use an external package (eg. mice) to generate imputed datasets, which can then be analyzed using the `semList` function. The `semTools` package contains several functions to do this automatically. Another option (with continuous data) is to use "two.stage" or "robust.two.stage". In this approach, we first estimate the sample statistics (mean vector, variance-covariance matrix) using an EM algorithm. Then, we use these estimated sample statistics as input for a regular analysis (as if the data were complete). The standard errors and test statistics are adjusted correctly to reflect the two-step procedure. The "robust.two.stage" option produces standard errors and a test statistic that are robust against non-normality. If (part of) the data is categorical, and the estimator is from the (W)LS family, the only option (besides listwise deletion) is "pairwise". In this three-step approach, missingness is only an issue in the first two steps. In the first step, we compute thresholds (for categorical variables) and means or intercepts (for continuous variables) using univariate information only. In this step, we simply ignore the missing values just like in `mean(x, na.rm = TRUE)`. In the second step, we compute polychoric/polyserial/pearson correlations using (only) two variables at a time. Here we use pairwise deletion: we only keep those observations for which both values are observed (not-missing). And this may change from pair to pair. By default, in the categorical case we use `conditional.x = TRUE`. Therefore, any cases with missing values on the exogenous covariates will be deleted listwise from the data first. Finally, if the estimator is "PML", the available options are "pairwise", "available.cases" and "doubly.robust". See the PML tutorial on the lavaan website for more information about these approaches.
- sampling.weights.normalization:** If "none", the sampling weights (if provided) will not be transformed. If "total", the sampling weights are normalized by dividing by the total sum of the weights, and multiplying again by the total sample size. If "group", the sampling weights are normalized per group: by dividing by the sum of the weights (in each group), and multiplying again by the group size. The default is "total".
- samplestats:** Logical. If FALSE, no sample statistics will be computed (and no estimation can take place). This can be useful when only a dummy lavaan object is requested, without any computations. The default is TRUE.

Data summary options:

sample.cov.rescale: If TRUE, the sample covariance matrix provided by the user is internally rescaled by multiplying it with a factor $(N-1)/N$. If "default", the value is set depending on the estimator and the likelihood option: it is set to TRUE if maximum likelihood estimation is used and likelihood="normal", and FALSE otherwise.

ridge: Logical. If TRUE a small constant value will be added the diagonal elements of the covariance (or correlation) matrix before analysis. The value can be set using the ridge.constant option.

ridge.constant: Numeric. Small constant used for ridgeing. The default value is 1e-05.

Multiple group options:

group.label: A character vector. The user can specify which group (or factor) levels need to be selected from the grouping variable, and in which order. If missing, all grouping levels are selected, in the order as they appear in the data.

group.equal: A vector of character strings. Only used in a multiple group analysis. Can be one or more of the following: "loadings", "composite.loadings", "intercepts", "means", "thresholds", "regressions", "residuals", "residual.covariances", "lv.variances" or "lv.covariances", specifying the pattern of equality constraints across multiple groups.

group.partial: A vector of character strings containing the labels of the parameters which should be free in all groups (thereby overriding the group.equal argument for some specific parameters).

group.w.free: Logical. If TRUE, the group frequencies are considered to be free parameters in the model. In this case, a Poisson model is fitted to estimate the group frequencies. If FALSE (the default), the group frequencies are fixed to their observed values.

Estimation options:

estimator: The estimator to be used. Can be one of the following: "ML" for maximum likelihood, "GLS" for (normal theory) generalized least squares, "WLS" for weighted least squares (sometimes called ADF estimation), "ULS" for unweighted least squares, "DWLS" for diagonally weighted least squares, and "DLS" for distributionally-weighted least squares. These are the main options that affect the estimation. For convenience, the "ML" option can be extended as "MLM", "MLMV", "MLMVS", "MLF", and "MLR". The estimation will still be plain "ML", but now with robust standard errors and a robust (scaled) test statistic. For "MLM", "MLMV", "MLMVS", classic robust standard errors are used (se="robust.sem"); for "MLF", standard errors are based on first-order derivatives (information = "first.order"); for "MLR", 'Huber-White' robust standard errors are used (se="robust.huber.white"). In addition, "MLM" will compute a Satorra-Bentler scaled (mean adjusted) test statistic (test="satorra.bentler"), "MLMVS" will compute a mean and variance adjusted test statistic (Satterthwaite style) (test="mean.var.adjusted"), "MLMV" will compute a mean and variance adjusted test statistic (scaled and shifted) (test="scaled.shifted"), and "MLR" will compute a test statistic which is asymptotically equivalent to the Yuan-Bentler T2-star test statistic (test="yuan.bentler.mplus"). Analogously, the estimators "WLSM" and "WLSMV" imply the "DWLS" estimator (not the "WLS" estimator) with robust standard errors and a mean or mean and variance adjusted test statistic. Estimators "ULSM" and "ULSMV" imply the "ULS" estimator with robust standard errors and a mean or mean and variance adjusted test statistic.

- likelihood:** Only relevant for ML estimation. If "wishart", the wishart likelihood approach is used. In this approach, the covariance matrix has been divided by N-1, and both standard errors and test statistics are based on N-1. If "normal", the normal likelihood approach is used. Here, the covariance matrix has been divided by N, and both standard errors and test statistics are based on N. If "default", it depends on the mimic option: if mimic="lavaan" or mimic="Mplus", normal likelihood is used; otherwise, wishart likelihood is used.
- link:** Not used yet. This is just a placeholder until the MML estimator is back.
- information:** If "expected", the expected information matrix is used (to compute the standard errors). If "observed", the observed information matrix is used. If "first.order", the information matrix is based on the outer product of the casewise scores. See also the options "h1.information" and "observed.information" for further control. If "default", the value is set depending on the estimator, the missing argument, and the mimic option. If the argument is a vector with two elements, the first element is used for the computation of the standard errors, while the second element is used for the (robust) test statistic.
- h1.information:** If "structured" (the default), the unrestricted (h1) information part of the (expected, first.order or observed if h1 is used) information matrix is based on the structured, or model-implied statistics (model-implied covariance matrix, model-implied mean vector, etc.). If "unstructured", the unrestricted (h1) information part is based on sample-based statistics (observed covariance matrix, observed mean vector, etc.) If the argument is a vector with two elements, the first element is used for the computation of the standard errors, while the second element is used for the (robust) test statistic.
- observed.information:** If "hessian", the observed information matrix is based on the hessian of the objective function. If "h1", an approximation is used that is based on the observed information matrix of the unrestricted (h1) model. If the argument is a vector with two elements, the first element is used for the computation of the standard errors, while the second element is used for the (robust) test statistic.
- se:** If "standard", conventional standard errors are computed based on inverting the (expected, observed or first.order) information matrix. If "robust.sem", conventional robust standard errors are computed. If "robust.huber.white", standard errors are computed based on the 'mlr' (aka pseudo ML, Huber-White) approach. If "robust", either "robust.sem" or "robust.huber.white" is used depending on the estimator, the mimic option, and whether the data are complete or not. If "boot" or "bootstrap", bootstrap standard errors are computed using standard bootstrapping (unless Bollen-Stine bootstrapping is requested for the test statistic; in this case bootstrap standard errors are computed using model-based bootstrapping). If "none", no standard errors are computed.
- test:** Character vector. See the documentation of the [lavTest](#) function for a full list. Multiple names of test statistics can be provided. If "default", the value depends on the values of other arguments. See also the [lavTest](#) function to extract (alternative) test statistics from a fitted lavaan object.
- scaled.test:** Character. Choose the test statistic that will be scaled (if a scaled test statistic is requested). The default is "standard", but it could also be (for example) "Browne.residual.nt".
- gamma.n.minus.one** Logical. If TRUE, we divide the Gamma matrix by N-1 (instead of the default N).
- gamma.unbiased** Logical. If TRUE, we compute an unbiased version for the Gamma matrix. Only available for single-level complete data and when conditional.x = FALSE and fixed.x = FALSE (for now).

bootstrap: Number of bootstrap draws, if bootstrapping is used.

do.fit: If FALSE, the model is not fit, and the current starting values of the model parameters are preserved.

Optimization options:

control: A list containing control parameters passed to the external optimizer. By default, lavaan uses "nlminb". See the manpage of `nlminb` for an overview of the control parameters. If another (external) optimizer is selected, see the manpage for that optimizer to see the possible control parameters.

optim.method: Character. The optimizer that should be used. For unconstrained optimization or models with only linear equality constraints (i.e., the model syntax does not include any "=", ">" or "<" operators), the available options are "nlminb" (the default), "BFGS", "L-BFGS-B". These are all quasi-newton methods. A basic implementation of Gauss-Newton is also available (`optim.method = "GN"`). The latter is the default when `estimator = "DLS"`. For constrained optimization, the only available option is "nlminb.constr", which uses an augmented Lagrangian minimization algorithm.

optim.force.converged: Logical. If TRUE, pretend the model has converged, no matter what.

optim.dx.tol Numeric. Tolerance used for checking if the elements of the (unscaled) gradient are all zero (in absolute value). The default value is 0.001.

optim.gn.tol.x: Numeric. Only used when `optim.method = "GN"`. Optimization stops when the root mean square of the difference between the old and new parameter values are smaller than this tolerance value. Default is $1e-05$ for DLS estimation and $1e-07$ otherwise.

optim.gn.iter.max: Integer. Only used when `optim.method = "GN"`. The maximum number of GN iterations. The default is 200.

bounds: Only used if `optim.method = "nlminb"`. If logical: FALSE implies no bounds are imposed on the parameters. If TRUE, this implies `bounds = "wide"`. If character, possible options are "none" (the default), "standard", "wide", "pos.var", "pos.ov.var", and "pos.lv.var". If `bounds = "pos.ov.var"`, the observed variances are forced to be nonnegative. If `bounds = "pos.lv.var"`, the latent variances are forced to be nonnegative. If `bounds = "pos.var"`, both observed and latent variances are forced to be nonnegative. If `bounds = "standard"`, lower and upper bounds are computed for observed and latent variances, and factor loadings. If `bounds = "wide"`, lower and upper bounds are computed for observed and latent variances, and factor loadings; but the range of the bounds is enlarged (allowing again for slightly negative variances).

optim.bounds: List. This can be used instead of the `bounds` argument to allow more control. Possible elements of the list are `lower`, `upper`, `lower.factor` and `upper.factor`. All of these accept a vector. The `lower` and `upper` elements indicate for which type of parameters bounds should be computed. Possible choices are "ov.var", "lv.var", "loadings" and "covariances". The `lower.factor` and `upper.factor` elements should have the same length as the `lower` and `upper` elements respectively. They indicate the factor by which the range of the bounds should be enlarged (for example, 1.1 or 1.2; the default is 1.0). Other elements are `min.reliability.marker` which sets the lower bound for the reliability of the marker indicator (if any) of each factor (default is 0.1). Finally, the `min.var.lv.endo` element indicates the lower bound of the variance of any endogenous latent variance (default is 0.0).

Parallelization options (currently only used for bootstrapping):

- parallel** The type of parallel operation to be used (if any). If missing, the default is "no".
- ncpus** Integer: number of processes to be used in parallel operation: typically one would chose this to the number of available CPUs. By default this is the number of cores (as detected by `parallel::detectCores()`) minus one.
- cl** An optional **parallel** or **snow** cluster for use if `parallel = "snow"`. If not supplied, a cluster on the local machine is created for the duration of the `bootstrapLavaan` or `bootstrapLRT` call.
- iseed** An integer to set the seed. Or NULL if no reproducible results are needed. This works for both serial (non-parallel) and parallel settings. Internally, `RNGkind()` is set to "L'Ecuyer-CMRG" if `parallel = "multicore"`. If `parallel = "snow"` (under windows), `parallel::clusterSetRNGStream()` is called which automatically switches to "L'Ecuyer-CMRG". When `iseed` is not NULL, `.Random.seed` (if it exists) in the global environment is left untouched.

Categorical estimation options:

- zero.add:** A numeric vector containing two values. These values affect the calculation of polychoric correlations when some frequencies in the bivariate table are zero. The first value only applies for 2x2 tables. The second value for larger tables. This value is added to the zero frequency in the bivariate table. If "default", the value is set depending on the "mimic" option. By default, lavaan uses `zero.add = c(0.5, 0, 0)`.
- zero.keep.margins:** Logical. This argument only affects the computation of polychoric correlations for 2x2 tables with an empty cell, and where a value is added to the empty cell. If TRUE, the other values of the frequency table are adjusted so that all margins are unaffected. If "default", the value is set depending on the "mimic". The default is TRUE.
- zero.cell.warn:** Logical. Only used if some observed endogenous variables are categorical. If TRUE, give a warning if one or more cells of a bivariate frequency table are empty.
- allow.empty.cell:** Logical. If TRUE, ignore situations where an ordinal variable has fewer categories than expected, or where a category is empty in a specific group.

Starting values options:

- start:** If it is a character string, the two options are currently "simple" and "Mplus". In the first case, all parameter values are set to zero, except the factor loadings and (residual) variances, which are set to one. When `start` is "Mplus", the factor loadings are estimated using the `fabin3` estimator (tsls) per factor. The residual variances of observed variables are set to half the observed variance, and all other (residual) variances are set to 0.05. The remaining parameters (regression coefficients, covariances) are set to zero. If `start` is a fitted object of class `lavaan`, the estimated values of the corresponding parameters will be extracted. If it is a parameter table, for example the output of the `parameterEstimates()` function, the values of the `est` or `start` or `ustart` column (whichever is found first) will be extracted.
- rstarts:** Integer. The number of refits that lavaan should try with random starting values. Random starting values are computed by drawing random numbers from a uniform distribution. Correlations are drawn from the interval [-0.5, +0.5] and then converted to covariances. Lower and upper bounds for (residual) variances are computed just like the standard bounds in bounded estimation. Random starting values are not computed for regression coefficients (which are always zero) and factor loadings of higher-order constructs (which are always unity). From all the runs that converged, the final solution is the one that resulted in the smallest value for the discrepancy function.

Check options:

- `check.start`: Logical. If TRUE, the starting values are checked for possibly inconsistent values (for example values implying correlations larger than one). If needed, a warning is given.
- `check.gradient`: Logical. If TRUE, and the model converged, a warning is given if the optimizer decided that a (local) solution has been found, while not all elements of the (unscaled) gradient (as seen by the optimizer) are (near) zero, as they should be (the tolerance used is 0.001).
- `check.post`: Logical. If TRUE, and the model converged, a check is performed after (post) fitting, to verify if the solution is admissible. This implies that all variances are non-negative, and all the model-implied covariance matrices are positive (semi-)definite. For the latter test, we tolerate a tiny negative eigenvalue that is smaller than $.Machine\$double.eps^{(3/4)}$, treating it as being zero.
- `check.vcov`: Logical. If TRUE, and the model converged, we check if the variance-covariance matrix of the free parameters is positive definite. We take into account possible equality and active inequality constraints. If needed, a warning is given.
- `check.lv.names`: Logical. If TRUE, and latent variables are defined in the model, lavaan will stop with an error message if a latent variable name also occurs in the data (implying it is also an observed variable).

Verbosity options:

- `verbose`: If TRUE, show what lavaan is doing. During estimation, the function value is printed out during each iteration.
- `warn`: If FALSE, suppress all lavaan-specific warning messages.
- `debug`: If TRUE, debugging information is printed out.

Miscellaneous:

- `model.type`: Set the model type: possible values are "cfa", "sem" or "growth". This may affect how starting values are computed, and may be used to alter the terminology used in the summary output, or the layout of path diagrams that are based on a fitted lavaan object.
- `mimic`: If "Mplus", an attempt is made to mimic the Mplus program. If "EQS", an attempt is made to mimic the EQS program. If "default", the value is (currently) set to "lavaan", which is very close to "Mplus".
- `representation`: If "LISREL" the classical LISREL matrix representation is used to represent the model (using the all-y variant). No other options are available (for now).
- `implied`: Logical. If TRUE, compute the model-implied statistics, and store them in the implied slot.
- `h1`: Logical. If TRUE, compute the unrestricted model and store the unrestricted summary statistics (and perhaps a loglikelihood) in the h1 slot.
- `baseline`: Logical. If TRUE, compute a baseline model (currently always the independence model, assuming all variables are uncorrelated) and store the results in the baseline slot.
- `baseline.conditional.x.free.slopes`: Logical. If TRUE, and `conditional.x = TRUE`, the (default) baseline model will allow the slopestructure to be unrestricted.
- `store.vcov` Logical. If TRUE, and `se=` is not set to "none", store the full variance-covariance matrix of the model parameters in the `vcov` slot of the fitted lavaan object.
- `parser` Character. If "new" (the default), the new parser is used to parse the model syntax. If "old", the original (pre 0.6-18) parser is used.

See Also[lavaan](#)**Examples**

```
lavOptions()
lavOptions("std.lv")
lavOptions(c("std.lv", "orthogonal"))
```

lavPredict

Predict the values of latent variables (and their indicators).

Description

The main purpose of the `lavPredict()` function is to compute (or ‘predict’) estimated values for the latent variables in the model (‘factor scores’). NOTE: the goal of this function is NOT to predict future values of dependent variables as in the regression framework! (For models with only continuous observed variables, the function `lavPredictY()` supports this.)

Usage

```
lavPredict(object, newdata = NULL, type = "lv", method = "EBM",
           transform = FALSE, se = "none", acov = "none",
           label = TRUE, fsm = FALSE, mdist = FALSE, rel = FALSE,
           append.data = FALSE, assemble = FALSE,
           level = 1L, optim.method = "bfgs", ETA = NULL,
           drop.list.single.group = TRUE)
```

Arguments

<code>object</code>	An object of class lavaan .
<code>newdata</code>	An optional <code>data.frame</code> , containing the same variables as the <code>data.frame</code> used when fitting the model in <code>object</code> .
<code>type</code>	A character string. If "lv", estimated values for the latent variables in the model are computed. If "ov", model predicted values for the indicators of the latent variables in the model are computed. If "yhat", the estimated value for the observed indicators, given user-specified values for the latent variables provided by de ETA argument. If "fy", densities (or probabilities) for each observed indicator, given user-specified values for the latent variables provided by de ETA argument.
<code>method</code>	A character string. In the linear case (when the indicators are continuous), the possible options are "regression" or "Bartlett". In the categorical case, the two options are "EBM" for the Empirical Bayes Modal approach, and "ML" for the maximum likelihood approach.

transform	Logical. If TRUE, transform the factor scores (per group) so that their mean and variance-covariance matrix matches the model-implied mean and variance-covariance matrix. This may be useful if the individual factor scores will be used in a follow-up (regression) analysis. Note: the standard errors (if requested) and the factor score matrix (if requested) are not transformed (yet).
se	Character. If "none", no standard errors are computed. If "standard", naive standard errors are computed (assuming the parameters of the measurement model are known). The standard errors are returned as an attribute. Currently only available for complete continuous data.
acov	Similar to the "se" argument, but optionally returns the full sampling covariance matrix of factor scores as an attribute. Currently only available for complete continuous data.
label	Logical. If TRUE, the columns in the output are labeled.
fsm	Logical. If TRUE, return the factor score matrix as an attribute. Only for numeric data.
mdist	Logical. If TRUE, the (squared) Mahalanobis distances of the factor scores (if type = "lv") or the casewise residuals (if type = "resid") are returned as an attribute.
rel	Logical. Only used if type = "lv". If TRUE, the factor reliabilities are returned as an attribute. (The squared values are often called the factor determinacies.)
append.data	Logical. Only used when type = "lv". If TRUE, the original data (or the data provided in the newdata argument) is appended to the factor scores.
assemble	Logical. If TRUE, the separate multiple groups are reassembled again to form a single data.frame with a group column, having the same dimensions as the original (or newdata) dataset.
level	Integer. Only used in a multilevel SEM. If level = 1, only factor scores for latent variable defined at the first (within) level are computed; if level = 2, only factor scores for latent variables defined at the second (between) level are computed.
optim.method	Character string. Only used in the categorical case. If "nlminb" (the default in 0.5), the "nlminb()" function is used for the optimization. If "bfgs" or "BFGS" (the default in 0.6), the "optim()" function is used with the BFGS method.
ETA	An optional matrix or list, containing latent variable values for each observation. Used for computations when type = "ov".
drop.list.single.group	Logical. If FALSE, the results are returned as a list, where each element corresponds to a group (even if there is only a single group). If TRUE, the list will be unlisted if there is only a single group.

Details

The `predict()` function calls the `lavPredict()` function with its default options.

If there are no latent variables in the model, `type = "ov"` will simply return the values of the observed variables. Note that this function can not be used to 'predict' values of dependent variables, given the values of independent values (in the regression sense). In other words, the structural component is completely ignored (for now).

See Also

[lavPredictY](#) to predict y-variables given x-variables.

Examples

```

data(HolzingerSwineford1939)

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

fit <- cfa(HS.model, data = HolzingerSwineford1939)
head(lavPredict(fit))
head(lavPredict(fit, type = "ov"))

## -----
## merge factor scores to original data.frame
## -----

idx <- lavInspect(fit, "case.idx")
fscores <- lavPredict(fit)
## loop over factors
for (fs in colnames(fscores)) {
  HolzingerSwineford1939[idx, fs] <- fscores[ , fs]
}
head(HolzingerSwineford1939)

## multigroup models return a list of factor scores (one per group)
data(HolzingerSwineford1939)
mgfit <- update(fit, group = "school", group.equal = c("loadings", "intercepts"))

idx <- lavInspect(mgfit, "case.idx") # list: 1 vector per group
fscores <- lavPredict(mgfit)         # list: 1 matrix per group
## loop over groups and factors
for (g in seq_along(fscores)) {
  for (fs in colnames(fscores[[g]])) {
    HolzingerSwineford1939[ idx[[g]], fs] <- fscores[[g]][ , fs]
  }
}
head(HolzingerSwineford1939)

## -----
## Use factor scores in susequent models
## -----

## see Examples in semTools package: ?plausibleValues

```

<code>lavPredictY</code>	<i>Predict the values of y-variables given the values of x-variables</i>
--------------------------	--

Description

This function can be used to predict the values of (observed) y-variables given the values of (observed) x-variables in a structural equation model.

Usage

```
lavPredictY(object, newdata = NULL,
            ynames = lavNames(object, "ov.y"),
            xnames = lavNames(object, "ov.x"),
            method = "conditional.mean",
            label = TRUE, assemble = TRUE,
            force.zero.mean = FALSE,
            lambda = 0)
```

Arguments

<code>object</code>	An object of class <code>lavaan</code> .
<code>newdata</code>	An optional data.frame, containing the same variables as the data.frame that was used when fitting the model in <code>object</code> . This data.frame should also include the y-variables (although their values will be ignored). Note that if no meanstructure was used in the original fit, we will use the saturated sample means of the original fit as substitutes for the model-implied means. Alternatively, refit the model using <code>meanstructure = TRUE</code> .
<code>ynames</code>	The names of the observed variables that should be treated as the y-variables. It is for these variables that the function will predict the (model-based) values for each observation. Can also be a list to allow for a separate set of variable names per group (or block).
<code>xnames</code>	The names of the observed variables that should be treated as the x-variables. Can also be a list to allow for a separate set of variable names per group (or block).
<code>method</code>	A character string. The only available option for now is "conditional.mean". See Details.
<code>label</code>	Logical. If TRUE, the columns of the output are labeled.
<code>assemble</code>	Logical. If TRUE, the predictions of the separate multiple groups in the output are reassembled again to form a single data.frame with a group column, having the same dimensions as the original (or newdata) dataset.
<code>force.zero.mean</code>	Logical. Only relevant if there is no mean structure. If TRUE, the (model-implied) mean vector is set to the zero vector. If FALSE, the (model-implied) mean vector is set to the (unrestricted) sample mean vector.
<code>lambda</code>	Numeric. A lambda regularization penalty term.

Details

This function can be used for (SEM-based) out-of-sample predictions of outcome (y) variables, given the values of predictor (x) variables. This is in contrast to the `lavPredict()` function which (historically) only ‘predicts’ the (factor) scores for latent variables, ignoring the structural part of the model.

When `method = "conditional.mean"`, predictions (for y given x) are based on the (joint y and x) model-implied variance-covariance (Sigma) matrix and mean vector (Mu), and the standard expression for the conditional mean of a multivariate normal distribution. Note that if the model is saturated (and hence `df = 0`), the SEM-based predictions are identical to ordinary least squares predictions.

Lambda is a regularization penalty term to improve prediction accuracy that can be determined using the `lavPredictY_cv` function.

References

de Rooij, M., Karch, J.D., Fokkema, M., Bakk, Z., Pratiwi, B.C, and Kelderman, H. (2022) SEM-Based Out-of-Sample Predictions, Structural Equation Modeling: A Multidisciplinary Journal. DOI:10.1080/10705511.2022.2061494

Molina, M. D., Molina, L., & Zappaterra, M. W. (2024). Aspects of Higher Consciousness: A Psychometric Validation and Analysis of a New Model of Mystical Experience. doi:10.31219/osf.io/cgb6e

See Also

[lavPredict](#) to compute scores for latent variables.

[lavPredictY_cv](#) to determine an optimal lambda to increase prediction accuracy.

Examples

```
model <- '
# latent variable definitions
ind60 =~ x1 + x2 + x3
dem60 =~ y1 + a*y2 + b*y3 + c*y4
dem65 =~ y5 + a*y6 + b*y7 + c*y8

# regressions
dem60 ~ ind60
dem65 ~ ind60 + dem60

# residual correlations
y1 ~~ y5
y2 ~~ y4 + y6
y3 ~~ y7
y4 ~~ y8
y6 ~~ y8
'

fit <- sem(model, data = PoliticalDemocracy)

lavPredictY(fit, ynames = c("y5", "y6", "y7", "y8"),
```



```
xnames = c("x1", "x2", "x3", "y1", "y2", "y3", "y4"))
```

lavPredictY_cv	<i>Determine an optimal lambda penalty value through cross-validation</i>
----------------	---

Description

This function can be used to determine an optimal lambda value for the `lavPredictY` function based on cross-validation.

Usage

```
lavPredictY_cv(object, data = NULL,
               xnames = lavNames(object, "ov.x"),
               ynames = lavNames(object, "ov.y"),
               n.folds = 10L,
               lambda.seq = seq(0, 1, 0.1))
```

Arguments

<code>object</code>	An object of class <code>lavaan</code> .
<code>data</code>	A <code>data.frame</code> , containing the same variables as the <code>data.frame</code> that was used when fitting the model in <code>object</code> .
<code>xnames</code>	The names of the observed variables that should be treated as the x-variables. Can also be a list to allow for a separate set of variable names per group (or block).
<code>ynames</code>	The names of the observed variables that should be treated as the y-variables. It is for these variables that the function will predict the (model-based) values for each observation. Can also be a list to allow for a separate set of variable names per group (or block).
<code>n.folds</code>	Integer. The number of folds to be used during cross-validation.
<code>lambda.seq</code>	An R <code>seq()</code> containing the range of lambda penalty values to be tested during cross-validation.

Details

This function is used to generate an optimal lambda value for `lavPredictY` predictions to improve prediction accuracy.

References

de Rooij, M., Karch, J.D., Fokkema, M., Bakk, Z., Pratiwi, B.C, and Kelderman, H. (2022) SEM-Based Out-of-Sample Predictions, *Structural Equation Modeling: A Multidisciplinary Journal*. DOI:10.1080/10705511.2022.2061494

Molina, M. D., Molina, L., & Zappaterra, M. W. (2024). Aspects of Higher Consciousness: A Psychometric Validation and Analysis of a New Model of Mystical Experience. doi:10.31219/osf.io/cgb6e

See Also

[lavPredictY](#) to predict the values of (observed) y-variables given the values of (observed) x-variables in a structural equation model.

Examples

```
colnames(PoliticalDemocracy) <- c("z1", "z2", "z3", "z4",
                                "y1", "y2", "y3", "y4",
                                "x1", "x2", "x3")

model <- '
# latent variable definitions
ind60 =~ x1 + x2 + x3
dem60 =~ z1 + z2 + z3 + z4
dem65 =~ y1 + y2 + y3 + y4
# regressions
dem60 ~ ind60
dem65 ~ ind60 + dem60
# residual correlations
z1 ~~ y1
z2 ~~ z4 + y2
z3 ~~ y3
z4 ~~ y4
y2 ~~ y4
'

fit <- sem(model, data = PoliticalDemocracy, meanstructure = TRUE)

percent <- 0.5
nobs <- lavInspect(fit, "ntotal")
idx <- sort(sample(x = nobs, size = floor(percent*nobs)))

xnames = c("z1", "z2", "z3", "z4", "x1", "x2", "x3")
ynames = c("y1", "y2", "y3", "y4")

reg.results <- lavPredictY_cv(
  fit,
  PoliticalDemocracy[-idx, ],
  xnames = xnames,
  ynames = ynames,
  n.folds = 10L,
  lambda.seq = seq(from = .6, to = 2.5, by = .1)
)
lam <- reg.results$lambda.min

lavPredictY(fit, newdata = PoliticalDemocracy[idx,],
            ynames = ynames,
            xnames = xnames,
            lambda = lam)
```

lavResiduals	<i>Residuals</i>
--------------	------------------

Description

'lavResiduals' provides model residuals and standardized residuals from a fitted lavaan object, as well as various summaries of these residuals.

The 'residuals()' (and 'resid()') methods are just shortcuts to this function with a limited set of arguments.

Usage

```
lavResiduals(object, type = "cor.bentler", custom.rmr = NULL,
             se = FALSE, zstat = TRUE, summary = TRUE, h1.acov = "unstructured",
             add.type = TRUE, add.labels = TRUE, add.class = TRUE,
             drop.list.single.group = TRUE,
             maximum.number = length(res.vech), output = "list")
```

Arguments

object	An object of class <code>lavaan</code> .
type	Character. If type = "raw", this function returns the raw (= unscaled) difference between the observed and the expected (model-implied) summary statistics, as well as the standardized version of these residuals. If type = "cor", or type = "cor.bollen", the observed and model implied covariance matrices are first transformed to a correlation matrix (using <code>cov2cor()</code>), before the residuals are computed. If type = "cor.bentler", both the observed and model implied covariance matrices are rescaled by dividing the elements by the square roots of the corresponding variances of the observed covariance matrix.
custom.rmr	list. Not used yet.
se	Logical. If TRUE, show the estimated standard errors for the residuals.
zstat	Logical. If TRUE, show the standardized residuals, which are the raw residuals divided by the corresponding (estimated) standard errors.
summary	Logical. If TRUE, show various summaries of the (possibly scaled) residuals. When type = "raw", we compute the RMR. When type = "cor.bentler", we compute the SRMR. When type = "cor.bollen", we compute the CRMR. An unbiased version of these summaries is also computed, as well as a standard error, a z-statistic and a p-value for the test of exact fit based on these summaries.
h1.acov	Character. If "unstructured", the observed summary statistics are used as consistent estimates of the corresponding (unrestricted) population statistics. If "structured", the model-implied summary statistics are used as consistent estimates of the corresponding (unrestricted) population statistics. This affects the way the asymptotic variance matrix of the summary statistics is computed.
add.type	Logical. If TRUE, show the type of residuals in the output.

<code>add.labels</code>	If TRUE, variable names are added to the vectors and/or matrices.
<code>add.class</code>	If TRUE, vectors are given the 'lavaan.vector' class; matrices are given the 'lavaan.matrix' class, and symmetric matrices are given the 'lavaan.matrix.symmetric' class. This only affects the way they are printed on the screen.
<code>drop.list.single.group</code>	If FALSE, the results are returned as a list, where each element corresponds to a group (even if there is only a single group). If TRUE, the list will be unlisted if there is only a single group.
<code>maximum.number</code>	Integer. Only used if <code>output = "table"</code> . Show only the first maximum.number rows of the data.frame.
<code>output</code>	Character. By default, <code>output = "list"</code> , and the output is a list of elements. If <code>output = "table"</code> , only the residuals of the variance-covariance matrix are shown in a data.frame, sorted from high (in absolute value) to low.

Value

If `drop.list.single.group = TRUE`, a list of (residualized) summary statistics, including type, standardized residuals, and summaries. If `drop.list.single.group = FALSE`, the list of summary statistics is nested within a list for each group.

References

Bentler, P.M. and Dijkstra, T. (1985). Efficient estimation via linearization in structural models. In Krishnaiah, P.R. (Ed.), *Multivariate analysis - VI*, (pp. 9–42). New York, NY: Elsevier.

Ogasawara, H. (2001). Standard errors of fit indices using residuals in structural equation modeling. *Psychometrika*, 66(3), 421–436. doi:10.1007/BF02294443

Maydeu-Olivares, A. (2017). Assessing the size of model misfit in structural equation models. *Psychometrika*, 82(3), 533–558. doi:10.1007/s11336-016-9552-7

Standardized Residuals in *Mplus*. Document retrieved from URL <http://www.statmodel.com/download/StandardizedResidual>

Examples

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

fit <- cfa(HS.model, data = HolzingerSwineford1939)
lavResiduals(fit)
```

lavTables

lavaan frequency tables

Description

Frequency tables for categorical variables and related statistics.

Usage

```
lavTables(object, dimension = 2L, type = "cells", categorical = NULL,
          group = NULL, statistic = "default", G2.min = 3, X2.min = 3,
          p.value = FALSE, output = "data.frame", patternAsString = TRUE)
```

Arguments

object	Either a data.frame, or an object of class <code>lavaan</code> .
dimension	Integer. If 0L, display all response patterns. If 1L, display one-dimensional (one-way) tables; if 2L, display two-dimensional (two-way or pairwise) tables. For the latter, we can change the information per row: if type = "cells", each row is a cell in a pairwise table; if type = "table", each row is a table.
type	If "cells", display information for each cell in the (one-way or two-way) table. If "table", display information per table. If "pattern", display response patterns (implying "dimension = 0L").
categorical	Only used if object is a data.frame. Specify variables that need to be treated as categorical.
group	Only used if object is a data.frame. Specify a grouping variable.
statistic	Either a character string, or a vector of character strings requesting one or more statistics for each cell, pattern or table. Always available are X2 and G2 for the Pearson and LRT based goodness-of-fit statistics. A distinction is made between the unrestricted and restricted model. The statistics based on the former have an extension *.un, as in X2.un and G2.un. If object is a data.frame, the unrestricted versions of the statistics are the only ones available. For one-way tables, additional statistics are the thresholds (th.un and th). For two-way tables and type = "table", the following statistics are available: X2, G2, cor (polychoric correlation), RMSEA and the corresponding unrestricted versions (X2.un etc). Additional statistics are G2.average, G2.nlarge and G2.plarge statistics based on the cell values G2: G2.average is the average of the G2 values in each cell of the two-way table; G2.nlarge is the number of cells with a G2 value larger than G2.min, and G2.plarge is the proportion of cells with a G2 value larger than G2.min. A similar set of statistics based on X2 is also available. If "default", the selection of statistics (if any) depends on the dim and type arguments, and if the object is a data.frame or a fitted lavaan object.
G2.min	Numeric. All cells with a G2 statistic larger than this number are considered 'large', as reflected in the (optional) "G2.plarge" and "G2.nlarge" columns.
X2.min	Numeric. All cells with a X2 statistic larger than this number are considered 'large', as reflected in the (optional) "X2.plarge" and "X2.nlarge" columns.
p.value	Logical. If "TRUE", p-values are computed for requested statistics (eg G2 or X2) if possible.
output	If "data.frame", the output is presented as a data.frame where each row is either a cell, a table, or a response pattern, depending on the "type" argument. If "table", the output is presented as a table (or matrix) or a list of tables. Only a single statistic can be shown in this case, and if the statistic is empty, the observed frequencies are shown.

patternAsString

Logical. Only used for response patterns (dimension = 0L). If "TRUE", response patterns are displayed as a compact string. If "FALSE", as many columns as observed variables are displayed.

Value

If output = "data.frame", the output is presented as a data.frame where each row is either a cell, a table, or a response pattern, depending on the "type" argument. If output = "table" (only for two-way tables), a list of tables (if type = "cells") where each list element corresponds to a pairwise table, or if type = "table", a single table (per group). In both cases, the table entries are determined by the (single) statistic argument.

References

Joreskog, K.G. & Moustaki, I. (2001). Factor analysis of ordinal variables: A comparison of three approaches. *Multivariate Behavioral Research*, 36, 347-387.

See Also

[varTable](#).

Examples

```
HS9 <- HolzingerSwineford1939[,c("x1", "x2", "x3", "x4", "x5",
                                "x6", "x7", "x8", "x9")]
HSbinary <- as.data.frame( lapply(HS9, cut, 2, labels=FALSE) )

# using the data only
lavTables(HSbinary, dim = 0L, categorical = names(HSbinary))
lavTables(HSbinary, dim = 1L, categorical = names(HSbinary), stat=c("th.un"))
lavTables(HSbinary, dim = 2L, categorical = names(HSbinary), type = "table")

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

fit <- cfa(HS.model, data=HSbinary, ordered=names(HSbinary))

lavTables(fit, 1L)
lavTables(fit, 2L, type="cells")
lavTables(fit, 2L, type="table", stat=c("cor.un", "G2", "cor"))
lavTables(fit, 2L, type="table", output="table", stat="X2")
```

lavTablesFitCp *Pairwise maximum likelihood fit statistics*

Description

Three measures of fit for the pairwise maximum likelihood estimation method that are based on likelihood ratios (LR) are defined: C_F , C_M , and C_P . Subscript F signifies a comparison of model-implied proportions of full response patterns with observed sample proportions, subscript M signifies a comparison of model-implied proportions of full response patterns with the proportions implied by the assumption of multivariate normality, and subscript P signifies a comparison of model-implied proportions of pairs of item responses with the observed proportions of pairs of item responses.

Usage

```
lavTablesFitCf(object)
lavTablesFitCp(object, alpha = 0.05)
lavTablesFitCm(object)
```

Arguments

object An object of class `lavaan`.
alpha The nominal level of significance of global fit.

Details

C_F : The C_F statistic compares the log-likelihood of the model-implied proportions (π_r) with the observed proportions (p_r) of the full multivariate responses patterns:

$$C_F = 2N \sum_r p_r \ln[p_r/\hat{\pi}_r],$$

which asymptotically has a chi-square distribution with

$$df_F = m^k - n - 1,$$

where k denotes the number of items with discrete response scales, m denotes the number of response options, and n denotes the number of parameters to be estimated. Notice that C_F results may be biased because of large numbers of empty cells in the multivariate contingency table.

C_M : The C_M statistic is based on the C_F statistic, and compares the proportions implied by the model of interest (Model 1) with proportions implied by the assumption of an underlying multivariate normal distribution (Model 0):

$$C_M = C_{F1} - C_{F0},$$

where C_{F0} is C_F for Model 0 and C_{F1} is C_F for Model 1. Statistic C_M has a chi-square distribution with degrees of freedom

$$df_M = k(k-1)/2 + k(m-1) - n_1,$$

where k denotes the number of items with discrete response scales, m denotes the number of response options, and $k(k-1)/2$ denotes the number of polychoric correlations, $k(m-1)$ denotes the number of thresholds, and n_1 is the number of parameters of the model of interest. Notice that C_M results may be biased because of large numbers of empty cells in the multivariate contingency table. However, bias may cancel out as both Model 1 and Model 0 contain the same pattern of empty responses.

C_P : With the C_P statistic we only consider pairs of responses, and compare observed sample proportions (p) with model-implied proportions of pairs of responses (π). For items i and j we obtain a pairwise likelihood ratio test statistic $C_{P_{ij}}$

$$C_{P_{ij}} = 2N \sum_{c_i=1}^m \sum_{c_j=1}^m p_{c_i, c_j} \ln[p_{c_i, c_j} / \hat{\pi}_{c_i, c_j}],$$

where m denotes the number of response options and N denotes sample size. The C_P statistic has an asymptotic chi-square distribution with degrees of freedom equal to the information ($m^2 - 1$) minus the number of parameters ($2(m-1)$ thresholds and 1 correlation),

$$df_P = m^2 - 2(m - 1) - 2.$$

As k denotes the number of items, there are $k(k-1)/2$ possible pairs of items. The C_P statistic should therefore be applied with a Bonferroni adjusted level of significance α^* , with

$$\alpha^* = \alpha / (k(k-1)/2),$$

to keep the family-wise error rate at α . The hypothesis of overall goodness-of-fit is tested at α and rejected as soon as C_P is significant at α^* for at least one pair of items. Notice that with dichotomous items, $m = 2$, and $df_P = 0$, so that hypothesis can not be tested.

References

- Barendse, M. T., Ligtvoet, R., Timmerman, M. E., & Oort, F. J. (2016). Structural Equation Modeling of Discrete data: Model Fit after Pairwise Maximum Likelihood. *Frontiers in psychology*, 7, 1-8.
- Joreskog, K. G., & Moustaki, I. (2001). Factor analysis of ordinal variables: A comparison of three approaches. *Multivariate Behavioral Research*, 36, 347-387.

See Also

[lavTables](#), [lavaan](#)

Examples

```
# Data
HS9 <- HolzingerSwineford1939[,c("x1", "x2", "x3", "x4", "x5",
                                "x6", "x7", "x8", "x9")]
HSbinary <- as.data.frame( lapply(HS9, cut, 2, labels=FALSE) )

# Single group example with one latent factor
HS.model <- ' trait =~ x1 + x2 + x3 + x4 '
```



```
fit <- cfa(HS.model, data=HSbinary[,1:4], ordered=names(HSbinary[,1:4]),
          estimator="PML")
lavTablesFitCm(fit)
lavTablesFitCp(fit)
lavTablesFitCf(fit)
```

lavTest	<i>Test of exact fit</i>
---------	--------------------------

Description

Compute a variety of test statistics evaluating the global fit of the model.

Usage

```
lavTest(lavobject, test = "standard", scaled.test = "standard",
        output = "list", drop.list.single = TRUE)
```

Arguments

lavobject	An object of class lavaan .
test	Character vector. Multiple names of test statistics can be provided. If "standard" is included, a conventional chi-square test is computed. If "Browne.residual.adf" is included, Browne's residual-based test statistic using ADF theory is computed. If "Browne.residual.nt" is included, Browne's residual-based test statistic using normal theory is computed. If "Satorra.Bentler" is included, a Satorra-Bentler scaled test statistic is computed. If "Yuan.Bentler" is included, a Yuan-Bentler scaled test statistic is computed. If "Yuan.Bentler.Mplus" is included, a test statistic is computed that is asymptotically equal to the Yuan-Bentler scaled test statistic. If "mean.var.adjusted" or "Satterthwaite" is included, a mean and variance adjusted test statistic is computed. If "scaled.shifted" is included, an alternative mean and variance adjusted test statistic is computed (as in Mplus version 6 or higher). If "boot" or "bootstrap" or "Bollen.Stine" is included, the Bollen-Stine bootstrap is used to compute the bootstrap probability value of the (regular) test statistic.
scaled.test	Character. Choose the test statistic that will be scaled (if a scaled test statistic is requested). The default is "standard", but it could also be (for example) "Browne.residual.nt".
output	Character. If "list" (the default), return a list with all test statistics. If "text", display the output as text with verbose descriptions (as in the summary output). If any scaled test statistics are included, they are printed first in a two-column format. Next come the other test statistics in a one-column format.
drop.list.single	Logical. Only used when output = "list". If TRUE and the list is of length one (i.e. only a single test statistic), drop the outer list. If FALSE, return a nested list with as many elements as we have test statistics.

Value

If output = "list": a nested list with test statistics, or if only a single test statistic is requested (and drop.list.single = TRUE), a list with details for this test statistic. If output = "text": the text is printed, and a nested list of test statistics (including an info attribute) is returned.

Examples

```
HS.model <- '
  visual =~ x1 + x2 + x3
  textual =~ x4 + x5 + x6
  speed  =~ x7 + x8 + x9
'
fit <- cfa(HS.model, data = HolzingerSwineford1939)
lavTest(fit, test = "browne.residual.adf")
```

lavTestLRT

LRT test

Description

LRT test for comparing (nested) lavaan models.

Usage

```
lavTestLRT(object, ..., method = "default", test = "default",
  A.method = "delta", scaled.shifted = TRUE,
  type = "Chisq", model.names = NULL)
anova(object, ...)
```

Arguments

object	An object of class <code>lavaan</code> .
...	additional objects of class <code>lavaan</code> .
method	Character string. The possible options are "satorra.bentler.2001", "satorra.bentler.2010", "satorra.2000", and "standard". See details.
test	Character string specifying which scaled test statistics to use, in case multiple scaled test= options were requested when fitting the model(s). See details.
A.method	Character string. The possible options are "exact" and "delta". This is only used when method = "satorra.2000". It determines how the Jacobian of the constraint function (the matrix A) will be computed. Note that if A.method = "exact", the models must be nested in the parameter sense, while if A.method = "delta", they only need to be nested in the covariance matrix sense.
scaled.shifted	Logical. Only used when method = "satorra.2000". If TRUE, we use a scaled and shifted test statistic; if FALSE, we use a mean and variance adjusted (Satterthwaite style) test statistic.

type	Character. If "Chisq", the test statistic for each model is the (scaled or unscaled) model fit test statistic. If "Cf", the test statistic for each model is computed by the <code>lavTablesFitCf</code> function. If "browne.residual.adf" (alias "browne") or "browne.residual.nt", the standard chi-squared difference is calculated from each model's residual-based statistic.
model.names	Character vector. If provided, use these model names in the first column of the anova table.

Details

The `anova` function for lavaan objects simply calls the `lavTestLRT` function, which has a few additional arguments.

The only `test=` options that currently have actual consequences are "satorra.bentler", "yuan.bentler", or "yuan.bentler.mplus" because "mean.var.adjusted" and "scaled.shifted" are currently distinguished by the `scaled.shifted` argument. See `lavOptions` for details about `test=` options implied by robust estimator= options). The "default" is to select the first available scaled statistic, if any. To check which test(s) were calculated when fitting your model(s), use `lavInspect(fit, "options")$test`.

If `type = "Chisq"` and the test statistics are scaled, a special scaled difference test statistic is computed. If `method` is "satorra.bentler.2001", a simple approximation is used described in Satorra & Bentler (2001). In some settings, this can lead to a negative test statistic. To ensure a positive test statistic, we can use the method proposed by Satorra & Bentler (2010). Alternatively, when `method="satorra.2000"`, the original formulas of Satorra (2000) are used. The latter is used for model comparison, when . . . contain additional (nested) models. Even when test statistics are scaled in object or . . . , users may request the `method="standard"` test statistic, without a robust adjustment.

Value

An object of class `anova`. When given a single argument, it simply returns the test statistic of this model. When given a sequence of objects, this function tests the models against one another, after reordering the models according to their degrees of freedom.

Note

If there is a `lavaan` model stored in `object@external$h1.model`, it will be added to . . .

References

- Satorra, A. (2000). Scaled and adjusted restricted tests in multi-sample analysis of moment structures. In Heijmans, R.D.H., Pollock, D.S.G. & Satorra, A. (eds.), *Innovations in multivariate statistical analysis: A Festschrift for Heinz Neudecker* (pp.233-247). London, UK: Kluwer Academic Publishers.
- Satorra, A., & Bentler, P. M. (2001). A scaled difference chi-square test statistic for moment structure analysis. *Psychometrika*, *66*(4), 507-514. doi:10.1007/BF02296192
- Satorra, A., & Bentler, P. M. (2010). Ensuring postiveness of the scaled difference chi-square test statistic. *Psychometrika*, *75*(2), 243-248. doi:10.1007/s113360099135y

Examples

```

HS.model <- '
  visual  =~ x1 + b1*x2 + x3
  textual =~ x4 + b2*x5 + x6
  speed   =~ x7 + b3*x8 + x9
,
fit1 <- cfa(HS.model, data = HolzingerSwineford1939)
fit0 <- cfa(HS.model, data = HolzingerSwineford1939,
            orthogonal = TRUE)
lavTestLRT(fit1, fit0)

## When multiple test statistics are selected when the model is fitted,
## use the type= and test= arguments to select a test for comparison.

## refit models, requesting 6 test statistics (in addition to "standard")
t6.1 <- cfa(HS.model, data = HolzingerSwineford1939,
            test = c("browne.residual.adf", "scaled.shifted", "mean.var.adjusted",
                    "satorra.bentler", "yuan.bentler", "yuan.bentler.mplus"))
t6.0 <- cfa(HS.model, data = HolzingerSwineford1939, orthogonal = TRUE,
            test = c("browne.residual.adf", "scaled.shifted", "mean.var.adjusted",
                    "satorra.bentler", "yuan.bentler", "yuan.bentler.mplus"))

## By default (test="default", type="Chisq"), the first scaled statistic
## requested will be used. Here, that is "scaled.shifted"
lavTestLRT(t6.1, t6.0)
## But even if "satorra.bentler" were requested first, method="satorra.2000"
## provides the scaled-shifted chi-squared difference test:
lavTestLRT(t6.1, t6.0, method = "satorra.2000")
## == lavTestLRT(update(t6.1, test = "scaled.shifted"), update(t6.0, test = "scaled.shifted"))

## The mean- and variance-adjusted (Satterthwaite) statistic implies
## scaled.shifted = FALSE
lavTestLRT(t6.1, t6.0, method = "satorra.2000", scaled.shifted = FALSE)

## Because "satorra.bentler" is not the first scaled test in the list,
## we MUST request it explicitly:
lavTestLRT(t6.1, t6.0, test = "satorra.bentler") # method="satorra.bentler.2001"
## == lavTestLRT(update(t6.1, test = "satorra.bentler"),
##               update(t6.0, test = "satorra.bentler"))
## The "strictly-positive test" is necessary when the above test is < 0:
lavTestLRT(t6.1, t6.0, test = "satorra.bentler", method = "satorra.bentler.2010")

## Likewise, other scaled statistics can be selected:
lavTestLRT(t6.1, t6.0, test = "yuan.bentler")
## == lavTestLRT(update(t6.1, test = "yuan.bentler"),
##               update(t6.0, test = "yuan.bentler"))
lavTestLRT(t6.1, t6.0, test = "yuan.bentler.mplus")
## == lavTestLRT(update(t6.1, test = "yuan.bentler.mplus"),
##               update(t6.0, test = "yuan.bentler.mplus"))

## To request the difference between Browne's (1984) residual-based statistics,

```

```
## rather than statistics based on the fitted model's discrepancy function,
## use the type= argument:
lavTestLRT(t6.1, t6.0, type = "browne.residual.adf")

## Despite requesting multiple robust tests, it is still possible to obtain
## the standard chi-squared difference test (i.e., without a robust correction)
lavTestLRT(t6.1, t6.0, method = "standard")
## == lavTestLRT(update(t6.1, test = "standard"), update(t6.0, test = "standard"))
```

lavTestScore	<i>Score test</i>
--------------	-------------------

Description

Score test (or Lagrange Multiplier test) for releasing one or more fixed or constrained parameters in model.

Usage

```
lavTestScore(object, add = NULL, release = NULL,
             univariate = TRUE, cumulative = FALSE,
             epc = FALSE, standardized = epc, cov.std = epc,
             verbose = FALSE, warn = TRUE, information = "expected")
```

Arguments

object	An object of class lavaan .
add	Either a character string (typically between single quotes) or a parameter table containing additional (currently fixed-to-zero) parameters for which the score test must be computed.
release	Vector of Integers. The indices of the constraints that should be released. The indices correspond to the order of the equality constraints as they appear in the parameter table.
univariate	Logical. If TRUE, compute the univariate score statistics, one for each constraints.
cumulative	Logical. If TRUE, order the univariate score statistics from large to small, and compute a series of multivariate score statistics, each time adding an additional constraint.
epc	Logical. If TRUE, and we are releasing existing constraints, compute the expected parameter changes for the existing (free) parameters, for each released constraint.
standardized	If TRUE, two extra columns (sepc.lv and sepc.all) in the \$epc table will contain standardized values for the EPCs. In the first column (sepc.lv), standardization is based on the variances of the (continuous) latent variables. In the second column (sepc.all), standardization is based on both the variances of both (continuous) observed and latent variables. (Residual) covariances are standardized using (residual) variances.

cov.std	Logical. See standardizedSolution .
verbose	Logical. Not used for now.
warn	Logical. If TRUE, print out warnings if they occur.
information	character indicating the type of information matrix to use (check lavInspect for available options). "expected" information is the default, which provides better control of Type I errors.

Details

This function can be used to compute both multivariate and univariate score tests. There are two modes: 1) releasing fixed-to-zero parameters (using the `add` argument), and 2) releasing existing equality constraints (using the `release` argument). The two modes can not be used simultaneously.

When adding new parameters, they should not already be part of the model (i.e. not listed in the parameter table). If you want to test for a parameter that was explicitly fixed to a constant (say to zero), it is better to label the parameter, and use an explicit equality constraint.

Value

A list containing at least one `data.frame`:

- `$test`: The total score test, with columns for the score test statistic (X^2), the degrees of freedom (df), and a p value under the χ^2 distribution (`p.value`).
- `$uni`: Optional (if `univariate=TRUE`). Each 1- df score test, equivalent to modification indices. If `epc=TRUE` when adding parameters (not when releasing constraints), an unstandardized EPC is provided for each added parameter, as would be returned by [modificationIndices](#).
- `$cumulative`: Optional (if `cumulative=TRUE`). Cumulative score tests.
- `$epc`: Optional (if `epc=TRUE`). Parameter estimates, expected parameter changes, and expected parameter values if all the tested constraints were freed.

References

Bentler, P. M., & Chou, C. P. (1993). Some new covariance structure model improvement statistics. Sage Focus Editions, 154, 235-255.

Examples

```
HS.model <- '
  visual =~ x1 + b1*x2 + x3
  textual =~ x4 + b2*x5 + x6
  speed  =~ x7 + b3*x8 + x9

  b1 == b2
  b2 == b3
'
fit <- cfa(HS.model, data=HolzingerSwineford1939)

# test 1: release both two equality constraints
lavTestScore(fit, cumulative = TRUE)
```

```

# test 2: the score test for adding two (currently fixed
# to zero) cross-loadings
newpar = '
  visual =~ x9
  textual =~ x3
  ,
lavTestScore(fit, add = newpar)

# equivalently, "add" can be a parameter table specifying parameters to free,
# but must include some additional information:
PT.add <- data.frame(lhs = c("visual", "textual"),
                    op = c("=~", "=~"),
                    rhs = c("x9", "x3"),
                    user = 10L, # needed to identify new parameters
                    free = 1, # arbitrary numbers > 0
                    start = 0) # null-hypothesized value

PT.add
lavTestScore(fit, add = PT.add) # same result as above

```

lavTestWald	<i>Wald test</i>
-------------	------------------

Description

Wald test for testing a linear hypothesis about the parameters of fitted lavaan object.

Usage

```
lavTestWald(object, constraints = NULL, verbose = FALSE)
```

Arguments

object	An object of class lavaan .
constraints	A character string (typically between single quotes) containing one or more equality constraints. See examples for more details.
verbose	Logical. If TRUE, print out the restriction matrix and the estimated restricted values.

Details

The constraints are specified using the "==" operator. Both the left-hand side and the right-hand side of the equality can contain a linear combination of model parameters, or a constant (like zero). The model parameters must be specified by their user-specified labels. Names of defined parameters (using the "!=" operator) can be included too.

Value

A list containing three elements: the Wald test statistic (stat), the degrees of freedom (df), and a p-value under the chi-square distribution (p.value).

Examples

```

HS.model <- '
  visual =~ x1 + b1*x2 + x3
  textual =~ x4 + b2*x5 + x6
  speed  =~ x7 + b3*x8 + x9
'

fit <- cfa(HS.model, data=HolzingerSwineford1939)

# test 1: test about a single parameter
# this is the 'chi-square' version of the
# z-test from the summary() output
lavTestWald(fit, constraints = "b1 == 0")

# test 2: several constraints
con = '
  2*b1 == b3
  b2 - b3 == 0
'

lavTestWald(fit, constraints = con)

```

lav_constraints

*Utility Functions: Constraints***Description**

Utility functions for equality and inequality constraints.

Usage

```

lav_constraints_parse(partable = NULL, constraints = NULL, theta = NULL,
  debug = FALSE)
lav_partable_constraints_ceq(partable, con = NULL, debug = FALSE,
  txtOnly = FALSE)
lav_partable_constraints_ciq(partable, con = NULL, debug = FALSE,
  txtOnly = FALSE)
lav_partable_constraints_def(partable, con = NULL, debug = FALSE,
  txtOnly = FALSE)

```

Arguments

partable	A lavaan parameter table.
constraints	A character string containing the constraints.
theta	A numeric vector. Optional vector with values for the model parameters in the parameter table.
debug	Logical. If TRUE, show debugging information.
con	An optional partable where the operator is one of '==', '>', '<' or ':='
txtOnly	Logical. If TRUE, only the body of the function is returned as a character string. If FALSE, a function is returned.

Details

This is a collection of lower-level constraint related functions that are used in the lavaan code. They are made public per request of package developers. Below is a brief description of what they do:

The `lav_constraints_parse` function parses the constraints specification (provided as a string, see example), and generates a list with useful information about the constraints.

The `lav_partable_constraints_ceq` function creates a function which takes the (unconstrained) parameter vector as input, and returns the slack values for each equality constraint. If the equality constraints hold perfectly, this function returns zeroes.

The `lav_partable_constraints_ciq` function creates a function which takes the (unconstrained) parameter vector as input, and returns the slack values for each inequality constraint.

The `lav_partable_constraints_def` function creates a function which takes the (unconstrained) parameter vector as input, and returns the computed values of the defined parameters.

Examples

```
myModel <- 'x1 ~ a*x2 + b*x3 + c*x4'
myParTable <- lavaanify(myModel, as.data.frame. = FALSE)
con <- ' a == 2*b
      b - c == 5 '
conInfo <- lav_constraints_parse(myParTable, constraints = con)

myModel2 <- 'x1 ~ a*x2 + b*x3 + c*x4
           a == 2*b
           b - c == 5 '
ceq <- lav_partable_constraints_ceq(partable = lavaanify(myModel2))
ceq( c(2,3,4) )
```

 lav_data

lavaan data functions

Description

Utility functions related to the Data slot

Usage

```
# update data slot with new data (of the same size)
lav_data_update(lavdata = NULL, newX = NULL, BOOT.idx = NULL, lavoptions = NULL)
```

Arguments

<code>lavdata</code>	A <code>lavdata</code> object.
<code>newX</code>	A list of (new) data matrices (per group) of the same size. They will replace the data stored in the internal <code>dataslot</code> .
<code>BOOT.idx</code>	A list of integers. If bootstrapping was used to produce the data in <code>newX</code> , use these indices to adapt the remaining slots.
<code>lavoptions</code>	A named list. The Options list from a lavaan object.

Examples

```
# generate syntax for an independence model
HS.model <- ' visual  =~ x1 + x2 + x3
            textual =~ x4 + x5 + x6
            speed   =~ x7 + x8 + x9 '

fit <- cfa(HS.model, data=HolzingerSwineford1939)

# extract data slot and options
lavdata <- fit@Data
lavoptions <- lavInspect(fit, "options")

# create bootstrap sample
boot.idx <- sample(x = nobs(fit), size = nobs(fit), replace = TRUE)
newX <- list(lavdata@X[[1]][boot.idx,])

# generate update lavdata object
newdata <- lav_data_update(lavdata = lavdata, newX = newX,
                          lavoptions = lavoptions)
```

lav_export_estimation *lav_export_estimation*

Description

lavaan provides a range of optimization methods with the `optim.method` argument (nlminb, BFGS, L-BFGS-B, GN, and nlminb.constr). ‘`lav_export_estimation`’ allows exporting objects and functions necessary to pass a lavaan model into any optimizer that takes a combination of (1) starting values, (2) fit-function, (3) gradient-function, and (4) upper and lower bounds. This allows testing new optimization frameworks.

Usage

```
lav_export_estimation(lavaan_model)
```

Arguments

```
lavaan_model    a fitted lavaan model
```

Value

List with:

- `get_coef` - When working with equality constraints, lavaan internally uses some transformations. `get_coef` is a functions that recreates the `coef` function for the parameters.
- `starting_values` - `starting_values` to be used in the optimization
- `objective_function` - objective function, expecting the current parameter values and the lavaan model

- gradient_function - gradient function, expecting the current parameter values and the lavaan model
- lower - lower bounds for parameters
- upper - upper bound for parameters

Examples

```

library(lavaan)
model <- '
  # latent variable definitions
  ind60 =~ x1 + x2 + x3
  dem60 =~ y1 + y2 + y3 + y4
  dem65 =~ y5 + a*y6 + y7 + y8

  # regressions
  dem60 ~ ind60
  dem65 ~ ind60 + dem60
'

fit <- sem(model,
           data = PoliticalDemocracy,
           do.fit = FALSE)

est <- lav_export_estimation(lavaan_model = fit)

# The starting values are:
est$starting_values
# Note that these do not have labels (and may also differ from coef(fit)
# in case of equality constraints):
coef(fit)
# To get the same parameters, use:
est$get_coef(parameter_values = est$starting_values,
             lavaan_model = fit)

# The objective function can be used to compute the fit at the current estimates:
est$objective_function(parameter_values = est$starting_values,
                      lavaan_model = fit)

# The gradient function can be used to compute the gradients at the current estimates:
est$gradient_function(parameter_values = est$starting_values,
                     lavaan_model = fit)

# Together, these elements provide the means to estimate the parameters with a large
# range of optimizers. For simplicity, here is an example using optim:
est_fit <- optim(par = est$starting_values,
               fn = est$objective_function,
               gr = est$gradient_function,
               lavaan_model = fit,
               method = "BFGS")
est$get_coef(parameter_values = est_fit$par,
             lavaan_model = fit)

```

```

# This is identical to
coef(sem(model,
        data = PoliticalDemocracy))

# Example using ridge regularization for parameter a
fn_ridge <- function(parameter_values, lavaan_model, est, lambda){
  return(est$objective_function(parameter_values = parameter_values,
                               lavaan_model = lavaan_model) + lambda * parameter_values[6]^2)
}
ridge_fit <- optim(par = est$get_coef(est$starting_values,
                                   lavaan_model = fit),
                 fn = fn_ridge,
                 lavaan_model = fit,
                 est = est,
                 lambda = 10)
est$get_coef(parameter_values = ridge_fit$par,
             lavaan_model = fit)

```

lav_func

Utility Functions: Gradient and Jacobian

Description

Utility functions for computing the gradient of a scalar-valued function or the Jacobian of a vector-valued function by numerical approximation.

Usage

```

lav_func_gradient_complex(func, x, h = .Machine$double.eps, ...,
                          fallback.simple = TRUE)
lav_func_jacobian_complex(func, x, h = .Machine$double.eps, ...,
                           fallback.simple = TRUE)

lav_func_gradient_simple(func, x, h = sqrt(.Machine$double.eps), ...)
lav_func_jacobian_simple(func, x, h = sqrt(.Machine$double.eps), ...)

```

Arguments

func	A real-valued function returning a numeric scalar or a numeric vector.
x	A numeric vector: the point(s) at which the gradient/Jacobian of the function should be computed.
h	Numeric value representing a small change in ‘x’ when computing the gradient/Jacobian.
...	Additional arguments to be passed to the function ‘func’.
fallback.simple	Logical. If TRUE, and the function evaluation fails, we call the corresponding simple (non-complex) method instead.

Details

The complex versions use complex numbers to gain more precision, while retaining the simplicity (and speed) of the simple forward method (see references). These functions were added to lavaan (around 2012) when the complex functionality was not part of the numDeriv package. They were used internally, and made public in 0.5-17 per request of other package developers.

References

Squire, W. and Trapp, G. (1998). Using Complex Variables to Estimate Derivatives of Real Functions. *SIAM Review*, 40(1), 110-112.

Examples

```
# very accurate complex method
lav_func_gradient_complex(func = exp, x = 1) - exp(1)

# less accurate forward method
lav_func_gradient_simple(func = exp, x = 1) - exp(1)

# very accurate complex method
diag(lav_func_jacobian_complex(func = exp, x = c(1,2,3))) - exp(c(1,2,3))

# less accurate forward method
diag(lav_func_jacobian_simple(func = exp, x = c(1,2,3))) - exp(c(1,2,3))
```

lav_matrix

Utility Functions: Matrices and Vectors

Description

Utility functions for Matrix and Vector operations.

Usage

```
# matrix to vector
lav_matrix_vec(A)
lav_matrix_vecr(A)
lav_matrix_vech(S, diagonal = TRUE)
lav_matrix_vechr(S, diagonal = TRUE)

# matrix/vector indices
lav_matrix_vech_idx(n = 1L, diagonal = TRUE)
lav_matrix_vech_row_idx(n = 1L, diagonal = TRUE)
lav_matrix_vech_col_idx(n = 1L, diagonal = TRUE)
lav_matrix_vechr_idx(n = 1L, diagonal = TRUE)
lav_matrix_vechr_u_idx(n = 1L, diagonal = TRUE)
lav_matrix_diag_idx(n = 1L)
lav_matrix_diagh_idx(n = 1L)
```

```

lav_matrix_antidiag_idx(n = 1L)

# vector to matrix
lav_matrix_vech_reverse(x, diagonal = TRUE)
lav_matrix_vechru_reverse(x, diagonal = TRUE)
lav_matrix_upper2full(x, diagonal = TRUE)
lav_matrix_vechr_reverse(x, diagonal = TRUE)
lav_matrix_vechu_reverse(x, diagonal = TRUE)
lav_matrix_lower2full(x, diagonal = TRUE)

# the duplication matrix
lav_matrix_duplication(n = 1L)
lav_matrix_duplication_pre(A = matrix(0,0,0))
lav_matrix_duplication_post(A = matrix(0,0,0))
lav_matrix_duplication_pre_post(A = matrix(0,0,0))
lav_matrix_duplication_ginv(n = 1L)
lav_matrix_duplication_ginv_pre(A = matrix(0,0,0))
lav_matrix_duplication_ginv_post(A = matrix(0,0,0))
lav_matrix_duplication_ginv_pre_post(A = matrix(0,0,0))

# the commutation matrix
lav_matrix_commutation(m = 1L, n = 1L)
lav_matrix_commutation_pre(A = matrix(0,0,0))
lav_matrix_commutation_post(A = matrix(0,0,0))
lav_matrix_commutation_pre_post(A = matrix(0,0,0))
lav_matrix_commutation_mn_pre(A, m = 1L, n = 1L)

# sample statistics
lav_matrix_cov(Y, Mu = NULL)

# other matrix operations
lav_matrix_symmetric_sqrt(S = matrix(0,0,0))
lav_matrix_orthogonal_complement(A = matrix(0,0,0))
lav_matrix_bdiag(...)
lav_matrix_trace(..., check = TRUE)

```

Arguments

A	A general matrix.
S	A symmetric matrix.
Y	A matrix representing a (numeric) dataset.
diagonal	Logical. If TRUE, include the diagonal.
n	Integer. When it is the only argument, the dimension of a square matrix. If m is also provided, the number of column of the matrix.
m	Integer. The number of rows of a matrix.
x	Numeric. A vector.

Mu	Numeric. If given, use Mu (instead of sample mean) to center, before taking the crossproduct.
...	One or more matrices, or a list of matrices.
check	Logical. If check = TRUE, we check if the (final) matrix is square.

Details

These are a collection of lower-level matrix/vector related functions that are used throughout the lavaan code. They are made public per request of package developers. Below is a brief description of what they do:

The `lav_matrix_vec` function implements the `vec` operator (for 'vectorization') and transforms a matrix into a vector by stacking the columns of the matrix one underneath the other.

The `lav_matrix_vecr` function is similar to the `lav_matrix_vec` function but transforms a matrix into a vector by stacking the rows of the matrix one underneath the other.

The `lav_matrix_vech` function implements the `vech` operator (for 'half vectorization') and transforms a symmetric matrix into a vector by stacking the columns of the matrix one underneath the other, but eliminating all supradiagonal elements. If `diagonal = FALSE`, the diagonal elements are also eliminated.

The `lav_matrix_vechr` function is similar to the `lav_matrix_vech` function but transforms a matrix into a vector by stacking the rows of the matrix one underneath the other, eliminating all supradiagonal elements.

The `lav_matrix_vech_idx` function returns the vector indices of the lower triangular elements of a symmetric matrix of size `n`, column by column.

The `lav_matrix_vech_row_idx` function returns the row indices of the lower triangular elements of a symmetric matrix of size `n`.

The `lav_matrix_vech_col_idx` function returns the column indices of the lower triangular elements of a symmetric matrix of size `n`.

The `lav_matrix_vechr_idx` function returns the vector indices of the lower triangular elements of a symmetric matrix of size `n`, row by row.

The `lav_matrix_vechu_idx` function returns the vector indices of the upper triangular elements of a symmetric matrix of size `n`, column by column.

The `lav_matrix_vechru_idx` function returns the vector indices of the upper triangular elements of a symmetric matrix of size `n`, row by row.

The `lav_matrix_diag_idx` function returns the vector indices of the diagonal elements of a symmetric matrix of size `n`.

The `lav_matrix_diagh_idx` function returns the vector indices of the lower part of a symmetric matrix of size `n`.

The `lav_matrix_antidiag_idx` function returns the vector indices of the anti diagonal elements a symmetric matrix of size `n`.

The `lav_matrix_vech_reverse` function (alias: `lav_matrix_vechru_reverse` and `lav_matrix_upper2full`) creates a symmetric matrix, given only upper triangular elements, row by row. If `diagonal = FALSE`, an diagonal with zero elements is added.

The `lav_matrix_vechr_reverse` (alias: `lav_matrix_vechu_reverse` and `lav_matrix_lower2full`) creates a symmetric matrix, given only the lower triangular elements, row by row. If `diagonal = FALSE`, an diagonal with zero elements is added.

The `lav_matrix_duplication` function generates the duplication matrix for a symmetric matrix of size n . This matrix duplicates the elements in `vech(S)` to create `vec(S)` (where S is symmetric). This matrix is very sparse, and should probably never be explicitly created. Use one of the functions below.

The `lav_matrix_duplication_pre` function computes the product of the transpose of the duplication matrix and a matrix A . The A matrix should have $n*n$ rows, where n is an integer. The duplication matrix is not explicitly created.

The `lav_matrix_duplication_post` function computes the product of a matrix A with the duplication matrix. The A matrix should have $n*n$ columns, where n is an integer. The duplication matrix is not explicitly created.

The `lav_matrix_duplication_pre_post` function first pre-multiplies a matrix A with the transpose of the duplication matrix, and then post multiplies the result again with the duplication matrix. A must be square matrix with $n*n$ rows and columns, where n is an integer. The duplication matrix is not explicitly created.

The `lav_matrix_duplication_ginv` function computes the generalized inverse of the duplication matrix. The matrix removes the duplicated elements in `vec(S)` to create `vech(S)`. This matrix is very sparse, and should probably never be explicitly created. Use one of the functions below.

The `lav_matrix_duplication_ginv_pre` function computes the product of the generalized inverse of the duplication matrix and a matrix A with $n*n$ rows, where n is an integer. The generalized inverse of the duplication matrix is not explicitly created.

The `lav_matrix_duplication_ginv_post` function computes the product of a matrix A (with $n*n$ columns, where n is an integer) and the transpose of the generalized inverse of the duplication matrix. The generalized inverse of the duplication matrix is not explicitly created.

The `lav_matrix_duplication_ginv_pre_post` function first pre-multiplies a matrix A with the transpose of the generalized inverse of the duplication matrix, and then post multiplies the result again with the transpose of the generalized inverse matrix. The matrix A must be square with $n*n$ rows and columns, where n is an integer. The generalized inverse of the duplication matrix is not explicitly created.

The `lav_matrix_commutation` function computes the commutation matrix which is a permutation matrix which transforms `vec(A)` (with m rows and n columns) into `vec(t(A))`.

The `lav_matrix_commutation_pre` function computes the product of the commutation matrix with a matrix A , without explicitly creating the commutation matrix. The matrix A must have $n*n$ rows, where n is an integer.

The `lav_matrix_commutation_post` function computes the product of a matrix A with the commutation matrix, without explicitly creating the commutation matrix. The matrix A must have $n*n$ rows, where n is an integer.

The `lav_matrix_commutation_pre_post` function first pre-multiplies a matrix A with the commutation matrix, and then post multiplies the result again with the commutation matrix, without explicitly creating the commutation matrix. The matrix A must have $n*n$ rows, where n is an integer.

The `lav_matrix_commutation_mn_pre` function computes the product of the commutation matrix with a matrix A, without explicitly creating the commutation matrix. The matrix A must have $m \times n$ rows, where m and n are integers.

The `lav_matrix_cov` function computes the sample covariance matrix of its input matrix, where the elements are divided by N (the number of rows).

The `lav_matrix_symmetric_sqrt` function computes the square root of a positive definite symmetric matrix (using an eigen decomposition). If some of the eigenvalues are negative, they are silently fixed to zero.

The `lav_matrix_orthogonal_complement` function computes an orthogonal complement of the matrix A, using a qr decomposition.

The `lav_matrix_bdiag` function constructs a block diagonal matrix from its arguments.

The `lav_matrix_trace` function computes the trace (the sum of the diagonal elements) of a single (square) matrix, or if multiple matrices are provided (either as a list, or as multiple arguments), we first compute their product (which must result in a square matrix), and then we compute the trace; if `check = TRUE`, we check if the (final) matrix is square.

References

Magnus, J. R. and H. Neudecker (1999). *Matrix Differential Calculus with Applications in Statistics and Econometrics*, Second Edition, John Wiley.

Examples

```
# upper elements of a 3 by 3 symmetric matrix (row by row)
x <- c(30, 16, 5, 10, 3, 1)
# construct full symmetric matrix
S <- lav_matrix_upper2full(x)

# compute the normal theory `Gamma' matrix given a covariance
# matrix (S), using the formula: Gamma = 2 * D^{+} (S %x% S) t(D^{+})
Gamma.NT <- 2 * lav_matrix_duplication_ginv_pre_post(S %x% S)
Gamma.NT
```

lav_model

lavaan model functions

Description

Utility functions related to internal model representation (lavmodel)

Usage

```
# set/get free parameters
lav_model_set_parameters(lavmodel, x = NULL)
lav_model_get_parameters(lavmodel, GLIST = NULL, type = "free",
                          extra = TRUE)
```

```
# compute model-implied statistics
lav_model_implied(lavmodel, GLIST = NULL, delta = TRUE)

# compute standard errors
lav_model_vcov_se(lavmodel, lavpartable, VCOV = NULL, BOOT = NULL)
```

Arguments

lavmodel	An internal representation of a lavaan model.
x	Numeric. A vector containing the values of all the free model parameters.
GLIST	List. A list of model matrices, similar to the output of <code>lavInspect(object, "est")</code> .
type	Character string. If "free", only return the free model parameters. If "user", return all the parameters (free and fixed) as they appear in the user-specified parameter table.
extra	Logical. If TRUE, also include values for rows in the parameter table where the operator is one of ":", "=", "<" or ">".
delta	Logical. If TRUE, and a Delta matrix is present in GLIST, use the (diagonal) values of the Delta matrix to rescale the covariance matrix. This is usually needed in the categorical setting to convert covariances to correlations.
lavpartable	A parameter table.
VCOV	Numeric matrix containing an estimate of the variance covariance matrix of the free model parameters.
BOOT	Numeric matrix containing the bootstrap based parameter estimates (in the columns) for each bootstrap sample (in the rows).

Examples

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

fit <- cfa(HS.model, data=HolzingerSwineford1939)
lavmodel <- fit@Model

est <- lav_model_get_parameters(lavmodel)
est
```

lav_partable *lavaan partable functions*

Description

Utility functions related to the parameter table (partable)

Usage

```

# extract information from a parameter table
lav_partable_df(partable)
lav_partable_ndat(partable)
lav_partable_npar(partable)
lav_partable_attributes(partable, pta = NULL)

# generate parameter labels
lav_partable_labels(partable, blocks = c("group", "level"),
                    group.equal = "", group.partial = "", type = "user")

# generate parameter table for specific models
lav_partable_independence(lavobject = NULL, lavdata = NULL,
                          lavpta = NULL, lavoptions = NULL, lavsamplestats = NULL,
                          lavh1 = NULL,
                          sample.cov = NULL, sample.mean = NULL, sample.slopes = NULL,
                          sample.th = NULL, sample.th.idx = NULL,
                          sample.cov.x = NULL, sample.mean.x = NULL)

lav_partable_unrestricted(lavobject = NULL, lavdata = NULL,
                          lavpta = NULL, lavoptions = NULL, lavsamplestats = NULL,
                          lavh1 = NULL,
                          sample.cov = NULL, sample.mean = NULL, sample.slopes = NULL,
                          sample.th = NULL, sample.th.idx = NULL,
                          sample.cov.x = NULL, sample.mean.x = NULL)

lav_partable_from_lm(object, est = FALSE, label = FALSE,
                    as.data.frame. = FALSE)

# complete a parameter table only containing a few columns (lhs,op,rhs)
lav_partable_complete(partable = NULL, start = TRUE)

# merge two parameter tables
lav_partable_merge(pt1 = NULL, pt2 = NULL, remove.duplicated = FALSE,
                  fromLast = FALSE, warn = TRUE)

# add a single parameter to an existing parameter table
lav_partable_add(partable = NULL, add = list())

```

Arguments

partable	A parameter table. See lavParTable for more information.
blocks	Character vector. Which columns in the parameter table should be taken to distinguish between different blocks of parameters (and hence be given different labels)? If "blocks" includes "group", a suffix ".g" and the group number (or group label) is added for the parameters of all but the first group. If "blocks" includes "level", a suffix ".l" and the level number is added for the parameters of all but the first level. If "blocks" includes, say "foo", a suffix ".foo" and

	the corresponding value of "foo" is added to all parameters.
group.equal	The same options can be used here as in the fitting functions. Parameters that are constrained to be equal across groups will be given the same label.
group.partial	A vector of character strings containing the labels of the parameters which should be free in all groups.
type	Character string. Can be either 'user' or 'free' to select all entries or only the free parameters from the parameter table respectively.
lavobject	An object of class 'lavaan'. If this argument is provided, it should be the only argument. All the values for the other arguments are extracted from this object.
lavdata	An object of class 'lavData'. The Data slot from a lavaan object.
lavoptions	A named list. The Options slot from a lavaan object.
lavsamplstats	An object of class 'lavSampleStats'. The SampleStats slot from a lavaan object.
lavh1	A named list. The h1 slot from a lavaan object.
lavpta	The pta (parameter table attributes) slot from a lavaan object.
sample.cov	Optional list of numeric matrices. Each list element contains a sample variance-covariance matrix for this group. If provided, these values will be used as starting values.
sample.mean	Optional list of numeric vectors. Each list element contains a sample mean vector for this group. If provided, these values will be used as starting values.
sample.slopes	Optional list of numeric matrices. Each list element contains the sample slopes for this group (only used when conditional.x = TRUE). If provided, these values will be used as starting values.
sample.th	Optional list of numeric vectors. Each list element contains a vector of sample thresholds for this group. If provided (and also sample.th.idx is provided), these values will be used as starting values.
sample.th.idx	Optional list of integers. Each list contains the threshold indices for this group.
sample.cov.x	Optional list of numeric matrices. Each list element contains a sample variance-covariance matrix for the exogenous variables for this group (only used when conditional.x = TRUE). If provided, these values will be used as starting values.
sample.mean.x	Optional list of numeric vectors. Each list element contains a sample mean vector for the exogenous variables for this group (only used when conditional.x = TRUE). If provided, these values will be used as starting values.
est	Logical. If TRUE, include the fitted estimates in the parameter table.
label	Logical. If TRUE, include parameter labels in the parameter table.
as.data.frame.	Logical. If TRUE, return the parameter table as a data.frame.
object	An object of class lm.
start	Logical. If TRUE, include a start column, based on the simple method for generating starting values.
pta	A list containing parameter attributes.
pt1	A parameter table.

pt2	A parameter table.
remove.duplicated	Logical. If TRUE, remove duplicated elements when merging two parameter tables.
fromLast	Logical. If TRUE, duplicated elements are considered from the bottom of the merged parameter table.
warn	Logical. If TRUE, a warning is produced when duplicated elements are found, when merging two parameter tables.
add	A named list. A single row of a parameter table as a named list.

Examples

```
# generate syntax for an independence model
HS.model <- ' visual =~ x1 + x2 + x3
            textual =~ x4 + x5 + x6
            speed  =~ x7 + x8 + x9 '

fit <- cfa(HS.model, data=HolzingerSwineford1939)
lav <- lav_partable_independence(fit)
as.data.frame(lav, stringsAsFactors = FALSE)

# how many free parameters?
lav_partable_npar(lav)

# how many sample statistics?
lav_partable_ndat(lav)
```

lav_samplestats *lavaan samplestats functions*

Description

Utility functions related to the sample statistics

Usage

```
# generate samplestats object from full data
lav_samplestats_from_data(lavdata = NULL, lavoptions = NULL,
                          WLS.V = NULL, NACOV = NULL)
```

Arguments

lavdata	A lavdata object.
lavoptions	A named list. The Options list from a lavaan object.
WLS.V	A user provided weight matrix.
NACOV	A user provided matrix containing the elements of (N times) the asymptotic variance-covariance matrix of the sample statistics. For a multiple group analysis, a list with an asymptotic variance-covariance matrix for each group.

Examples

```
# generate syntax for an independence model
HS.model <- ' visual  =~ x1 + x2 + x3
            textual =~ x4 + x5 + x6
            speed   =~ x7 + x8 + x9 '

fit <- cfa(HS.model, data=HolzingerSwineford1939)

# extract data slot and options
lavdata <- fit@Data
lavoptions <- lavInspect(fit, "options")

# generate sample statistics object
sampleStats <- lav_samplestats_from_data(lavdata = lavdata,
                                         lavoptions = lavoptions)
```

model.syntax

The Lavaan Model Syntax

Description

The lavaan model syntax describes a latent variable model. The function `lavaanify` turns it into a table that represents the full model as specified by the user. We refer to this table as the parameter table.

Usage

```
lavaanify(model = NULL, meanstructure = FALSE, int.ov.free = FALSE,
           int.lv.free = FALSE, marker.int.zero = FALSE,
           orthogonal = FALSE, orthogonal.y = FALSE,
           orthogonal.x = FALSE, orthogonal.efa = FALSE, std.lv = FALSE,
           correlation = FALSE, effect.coding = "", conditional.x = FALSE,
           fixed.x = FALSE, parameterization = "delta", constraints = NULL,
           ceq.simple = FALSE, auto = FALSE, model.type = "sem",
           auto.fix.first = FALSE, auto.fix.single = FALSE, auto.var = FALSE,
           auto.cov.lv.x = FALSE, auto.cov.y = FALSE, auto.th = FALSE,
           auto.delta = FALSE, auto.efa = FALSE,
           varTable = NULL, ngroups = 1L, nthresholds = NULL,
           group.equal = NULL, group.partial = NULL, group.w.free = FALSE,
           debug = FALSE, warn = TRUE, as.data.frame. = TRUE)

lavParTable(model = NULL, meanstructure = FALSE, int.ov.free = FALSE,
            int.lv.free = FALSE, marker.int.zero = FALSE,
            orthogonal = FALSE, orthogonal.y = FALSE,
            orthogonal.x = FALSE, orthogonal.efa = FALSE, std.lv = FALSE,
            correlation = FALSE, effect.coding = "", conditional.x = FALSE,
            fixed.x = FALSE, parameterization = "delta", constraints = NULL,
            ceq.simple = FALSE, auto = FALSE, model.type = "sem",
```

```

auto.fix.first = FALSE, auto.fix.single = FALSE, auto.var = FALSE,
auto.cov.lv.x = FALSE, auto.cov.y = FALSE, auto.th = FALSE,
auto.delta = FALSE, auto.efa = FALSE,
varTable = NULL, ngroups = 1L, nthresholds = NULL,
group.equal = NULL, group.partial = NULL, group.w.free = FALSE,
debug = FALSE, warn = TRUE, as.data.frame. = TRUE)

```

```

lavParseModelString(model.syntax = '', as.data.frame. = FALSE,
                    parser = "new", warn = TRUE, debug = FALSE)

```

Arguments

<code>model</code>	A description of the user-specified model. Typically, the model is described using the lavaan model syntax; see details for more information. Alternatively, a parameter table (e.g., the output of <code>lavParseModelString</code> is also accepted.
<code>model.syntax</code>	The model syntax specifying the model. Must be a literal string.
<code>meanstructure</code>	If TRUE, intercepts/means will be added to the model both for both observed and latent variables.
<code>int.ov.free</code>	If FALSE, the intercepts of the observed variables are fixed to zero.
<code>int.lv.free</code>	If FALSE, the intercepts of the latent variables are fixed to zero.
<code>marker.int.zero</code>	Logical. Only relevant if the metric of each latent variable is set by fixing the first factor loading to unity. If TRUE, it implies <code>meanstructure = TRUE</code> and <code>std.lv = FALSE</code> , and it fixes the intercepts of the marker indicators to zero, while freeing the means/intercepts of the latent variables. Only works correctly for single group, single level models.
<code>orthogonal</code>	If TRUE, all covariances among latent variables are set to zero.
<code>orthogonal.y</code>	If TRUE, all covariances among endogenous latent variables only are set to zero.
<code>orthogonal.x</code>	If TRUE, all covariances among exogenous latent variables only are set to zero.
<code>orthogonal.efa</code>	If TRUE, all covariances among latent variables involved in rotation only are set to zero.
<code>std.lv</code>	If TRUE, the metric of each latent variable is determined by fixing their variances to 1.0. If FALSE, the metric of each latent variable is determined by fixing the factor loading of the first indicator to 1.0. If there are multiple groups, <code>std.lv = TRUE</code> and "loadings" is included in the <code>group.label</code> argument, then only the latent variances <i>i</i> of the first group will be fixed to 1.0, while the latent variances of other groups are set free.
<code>correlation</code>	If TRUE, a correlation structure is fitted. For continuous data, this implies that the (residual) variances are no longer parameters of the model.
<code>effect.coding</code>	Can be logical or character string. If logical and TRUE, this implies <code>effect.coding = c("loadings", "intercepts")</code> . If logical and FALSE, it is set equal to the empty string. If "loadings" is included, equality constraints are used so that the average of the factor loadings (per latent variable) equals 1. Note that this should not be used together with <code>std.lv = TRUE</code> . If "intercepts" is included,

equality constraints are used so that the sum of the intercepts (belonging to the indicators of a single latent variable) equals zero. As a result, the latent mean will be freely estimated and usually equal the average of the means of the involved indicators.

conditional.x	If TRUE, we set up the model conditional on the exogenous 'x' covariates; the model-implied sample statistics only include the non-x variables. If FALSE, the exogenous 'x' variables are modeled jointly with the other variables, and the model-implied statistics reflect both sets of variables.
fixed.x	If TRUE, the exogenous 'x' covariates are considered fixed variables and the means, variances and covariances of these variables are fixed to their sample values. If FALSE, they are considered random, and the means, variances and covariances are free parameters.
parameterization	Currently only used if data is categorical. If "delta", the delta parameterization is used. If "theta", the theta parameterization is used.
constraints	Additional (in)equality constraints. See details for more information.
ceq.simple	If TRUE, and no other general constraints are used in the model, simple equality constraints are represented in the parameter table as duplicated free parameters (instead of extra rows with op = "==").
auto	If TRUE, the default values are used for the auto.* arguments, depending on the value of model.type.
model.type	Either "sem" or "growth"; only used if auto=TRUE.
auto.fix.first	If TRUE, the factor loading of the first indicator is set to 1.0 for every latent variable.
auto.fix.single	If TRUE, the residual variance (if included) of an observed indicator is set to zero if it is the only indicator of a latent variable.
auto.var	If TRUE, the (residual) variances of both observed and latent variables are set free.
auto.cov.lv.x	If TRUE, the covariances of exogenous latent variables are included in the model and set free.
auto.cov.y	If TRUE, the covariances of dependent variables (both observed and latent) are included in the model and set free.
auto.th	If TRUE, thresholds for limited dependent variables are included in the model and set free.
auto.delta	If TRUE, response scaling parameters for limited dependent variables are included in the model and set free.
auto.efa	If TRUE, the necessary constraints are imposed to make the (unrotated) exploratory factor analysis blocks identifiable: for each block, factor variances are set to 1, factor covariances are constrained to be zero, and factor loadings are constrained to follow an echelon pattern.
varTable	The variable table containing information about the observed variables in the model.
ngroups	The number of (independent) groups.

nthresholds	Either a single integer or a named vector of integers. If nthresholds is a single integer, all endogenous variables are assumed to be ordered with nthresholds indicating the number of thresholds needed in the model. If nthresholds is a named vector, it indicates the number of thresholds for these ordered variables only. This argument should not be used in combination with varTable.
group.equal	A vector of character strings. Only used in a multiple group analysis. Can be one or more of the following: "loadings", "intercepts", "means", "regressions", "residuals" or "covariances", specifying the pattern of equality constraints across multiple groups. When (in the model syntax) a vector of labels is used as a modifier for a certain parameter, this will override the group.equal setting if it applies to this parameter. See also the Multiple groups section below for using modifiers in multiple groups.
group.partial	A vector of character strings containing the labels of the parameters which should be free in all groups (thereby overriding the group.equal argument for some specific parameters).
group.w.free	Logical. If TRUE, the group frequencies are considered to be free parameters in the model. In this case, a Poisson model is fitted to estimate the group frequencies. If FALSE (the default), the group frequencies are fixed to their observed values.
as.data.frame.	If TRUE, return the list of model parameters as a data.frame.
parser	Character. If "old", use the original/classic parser. If "new", use the new/ldw parser. The default is "new".
warn	If TRUE, some (possibly harmless) warnings are printed out.
debug	If TRUE, debugging information is printed out.

Details

The model syntax consists of one or more formula-like expressions, each one describing a specific part of the model. The model syntax can be read from a file (using [readLines](#)), or can be specified as a literal string enclosed by single quotes as in the example below.

```
myModel <- '
# 1. latent variable definitions
  f1 =~ y1 + y2 + y3
  f2 =~ y4 + y5 + y6
  f3 =~ y7 + y8 +
      y9 + y10
  f4 =~ y11 + y12 + y13

! this is also a comment

# 2. regressions
  f1 ~ f3 + f4
  f2 ~ f4
  y1 + y2 ~ x1 + x2 + x3

# 3. (co)variances
```

```

y1 ~~ y1
y2 ~~ y4 + y5
f1 ~~ f2

# 4. intercepts
f1 ~ 1; y5 ~ 1

# 5. thresholds
y11 | t1 + t2 + t3
y12 | t1
y13 | t1 + t2

# 6. scaling factors
y11 ~*~ y11
y12 ~*~ y12
y13 ~*~ y13

# 7. formative factors
f5 <~ z1 + z2 + z3 + z4

```

Blank lines and comments can be used in between the formulas, and formulas can be split over multiple lines. Both the sharp (#) and the exclamation (!) characters can be used to start a comment. Multiple formulas can be placed on a single line if they are separated by a semicolon (;).

There can be seven types of formula-like expressions in the model syntax:

1. Latent variable definitions: The "`~~`" operator can be used to define (continuous) latent variables. The name of the latent variable is on the left of the "`~~`" operator, while the terms on the right, separated by "`+`" operators, are the indicators of the latent variable. The operator "`~~`" can be read as "is manifested by".
2. Regressions: The "`~`" operator specifies a regression. The dependent variable is on the left of a "`~`" operator and the independent variables, separated by "`+`" operators, are on the right. These regression formulas are similar to the way ordinary linear regression formulas are used in R, but they may include latent variables. Interaction terms are currently not supported.
3. Variance-covariances: The "`~~~`" ('double tilde') operator specifies (residual) variances of an observed or latent variable, or a set of covariances between one variable, and several other variables (either observed or latent). Several variables, separated by "`+`" operators can appear on the right. This way, several pairwise (co)variances involving the same left-hand variable can be expressed in a single expression. The distinction between variances and residual variances is made automatically.
4. Intercepts: A special case of a regression formula can be used to specify an intercept (or a mean) of either an observed or a latent variable. The variable name is on the left of a "`~`" operator. On the right is only the number "1" representing the intercept. Including an intercept formula in the model automatically implies `meanstructure = TRUE`. The distinction between intercepts and means is made automatically.
5. Thresholds: The "`|`" operator can be used to define the thresholds of categorical endogenous variables (on the left hand side of the operator). By convention, the thresholds (on the right hand sided, separated by the "`+`" operator, are named "`t1`", "`t2`", etcetera.

6. Scaling factors: The " $\sim*$ " operator defines a scale factor. The variable name on the left hand side must be the same as the variable name on the right hand side. Scale factors are used in the Delta parameterization, in a multiple group analysis when factor indicators are categorical.
7. Formative factors: The " $\sim<$ " operator can be used to define a formative factor (on the right hand side of the operator), in a similar way to how a reflexive factor is defined (using the " $\sim=$ " operator). This is just syntax sugar to define a phantom latent variable (equivalent to using " $f \sim 0$ "). And in addition, the (residual) variance of the formative factor is fixed to zero.

There are 4 additional operators, also with left- and right-hand sides, that can be included in model syntax. Three of them are used to specify (in)equality constraints on estimated parameters ($\sim=$, $\sim>$, and $\sim<$), and those are demonstrated in a later section about **(In)equality constraints**. The final additional operator ($\sim:=$) can be used to define "new" parameters that are functions of one or more other estimated parameters. The $\sim:=$ operator is demonstrated in a section about **User-defined parameters**.

Usually, only a single variable name appears on the left side of an operator. However, if multiple variable names are specified, separated by the "+" operator, the formula is repeated for each element on the left side (as for example in the third regression formula in the example above). The only exception are scaling factors, where only a single element is allowed on the left hand side.

In the right-hand side of these formula-like expressions, each element can be modified (using the " $\sim*$ " operator) by either a numeric constant, an expression resulting in a numeric constant, an expression resulting in a character vector, or one of three special functions: `start()`, `label()` and `equal()`. This provides the user with a mechanism to fix parameters, to provide alternative starting values, to label the parameters, and to define equality constraints among model parameters. All " $\sim*$ " expressions are referred to as *modifiers*. They are explained in more detail in the following sections.

Fixing parameters

It is often desirable to fix a model parameter that is otherwise (by default) free. Any parameter in a model can be fixed by using a modifier resulting in a numerical constraint. Here are some examples:

- Fixing the regression coefficient of the predictor x2:

$$y \sim x1 + 2.4*x2 + x3$$

- Specifying an orthogonal (zero) covariance between two latent variables:

$$f1 \sim\sim 0*f2$$

- Specifying an intercept and a linear slope in a growth model:

$$\begin{aligned} i &\sim 1*y11 + 1*y12 + 1*y13 + 1*y14 \\ s &\sim 0*y11 + 1*y12 + 2*y13 + 3*y14 \end{aligned}$$

Instead of a numeric constant, one can use a mathematical function that returns a numeric constant, for example `sqrt(10)`. Multiplying with NA will force the corresponding parameter to be free.

Additionally, the $\sim=$ operator can be used to set a *labeled* parameter equal to a specific numeric value. This will be demonstrated in the section below about **(In)equality constraints**.

Starting values

User-provided starting values can be given by using the special function `start()`, containing a numeric constant. For example:

```
y ~ x1 + start(1.0)*x2 + x3
```

Note that if a starting value is provided, the parameter is not automatically considered to be free.

Parameter labels and equality constraints

Each free parameter in a model is automatically given a name (or label). The name given to a model parameter consists of three parts, coerced to a single character vector. The first part is the name of the variable in the left-hand side of the formula where the parameter was implied. The middle part is based on the special ‘operator’ used in the formula. This can be either one of “= \sim ”, “ \sim ” or “ $\sim\sim$ ”. The third part is the name of the variable in the right-hand side of the formula where the parameter was implied, or “1” if it is an intercept. The three parts are pasted together in a single string. For example, the name of the fixed regression coefficient in the regression formula $y \sim x1 + 2.4*x2 + x3$ is the string “y~x2”. The name of the parameter corresponding to the covariance between two latent variables in the formula $f1 \sim\sim f2$ is the string “f1~f2”.

Although this automatic labeling of parameters is convenient, the user may specify its own labels for specific parameters simply by pre-multiplying the corresponding term (on the right hand side of the operator only) by a character string (starting with a letter). For example, in the formula $f1 \sim x1 + x2 + mylabel*x3$, the parameter corresponding with the factor loading of $x3$ will be named “mylabel”. An alternative way to specify the label is as follows: $f1 \sim x1 + x2 + label("mylabel")*x3$, where the label is the argument of special function `label()`; this can be useful if the label contains a space, or an operator (like “ \sim ”).

To constrain a parameter to be equal to another target parameter, there are two ways. If you have specified your own labels, you can use the fact that *equal labels imply equal parameter values*. If you rely on automatic parameter labels, you can use the special function `equal()`. The argument of `equal()` is the (automatic or user-specified) name of the target parameter. For example, in the confirmatory factor analysis example below, the intercepts of the three indicators of each latent variable are constrained to be equal to each other. For the first three, we have used the default names. For the last three, we have provided a custom label for the $y2a$ intercept.

```
model <- '
  # two latent variables with fixed loadings
  f1 =~ 1*y1a + 1*y1b + 1*y1c
  f2 =~ 1*y2a + 1*y2b + 1*y2c

  # intercepts constrained to be equal
  # using the default names
  y1a ~ 1
  y1b ~ equal("y1a~1") * 1
  y1c ~ equal("y1a~1") * 1

  # intercepts constrained to be equal
  # using a custom label
  y2a ~ int2*1
```

```

y2b ~ int2*1
y2c ~ int2*1

```

Multiple groups

In a multiple group analysis, modifiers that contain a single element should be replaced by a vector, having the same length as the number of groups. If you provide a single element, it will be recycled for all the groups. This may be dangerous, in particular when the modifier is a label. In that case, the (same) label is copied across all groups, and this would imply an equality constraint across groups. Therefore, when using modifiers in a multiple group setting, it is always safer (and cleaner) to specify the same number of elements as the number of groups. Consider this example with two groups:

```

HS.model <- ' visual  =~ x1 + 0.5*x2 + c(0.6, 0.8)*x3
             textual =~ x4 + start(c(1.2, 0.6))*x5 + x6
             speed   =~ x7 + x8 + c(x9.group1, x9.group2)*x9 '

```

In this example, the factor loading of the 'x2' indicator is fixed to the value 0.5 for both groups. However, the factor loadings of the 'x3' indicator are fixed to 0.6 and 0.8 for group 1 and group 2 respectively. The same logic is used for all modifiers. Note that character vectors can contain unquoted strings.

Multiple modifiers

In the model syntax, you can specify a variable more than once on the right hand side of an operator; therefore, several 'modifiers' can be applied simultaneously; for example, if you want to fix the value of a parameter and also label that parameter, you can use something like:

```
f1 =~ x1 + x2 + 4*x3 + x3.loading*x3
```

(In)equality constraints

The == operator can be used either to fix a parameter to a specific value, or to set an estimated parameter equal to another parameter. Adapting the example in the **Parameter labels and equality constraints** section, we could have used different labels for the second factor's intercepts:

```

y2a ~ int1*1
y2b ~ int2*1
y2c ~ int3*1

```

Then, we could fix the first intercept to zero by including in the syntax an operation that indicates the parameter's label equals that value:

```
int1 == 0
```

Whereas we could still estimate the other two intercepts under an equality constraint by setting their different labels equal to each other:

```
int2 == int3
```

Optimization can be less efficient when constraining parameters this way (see the documentation linked under **See also** for more information). But the flexibility might be advantageous. For example, the constraints could be specified in a separate character-string object, which can be passed to the `lavaan(..., constraints=)` argument, enabling users to compare results with(out) the constraints.

Inequality constraints work much the same way, using the `<` or `>` operator indicate which estimated parameter is hypothesized to be greater/less than either a specific value or another estimated parameter. For example, a variance can be constrained to be nonnegative:

```
y1a ~~ var1a*y1a
## hypothesized constraint:
var1a > 0
```

Or the factor loading of a particular indicator might be expected to exceed other indicators' loadings:

```
f1 =~ L1*y1a + L2*y1b + L3*y1c
## hypothesized constraints:
L1 > L2
L3 < L1
```

User-defined parameters

Functions of parameters can be useful to test particular hypotheses. Following from the Multiple groups example, we might be interested in which group's factor loading is larger (i.e., an estimate of differential item functioning (DIF) when the latent scales are linked by anchor items with equal loadings).

```
speed =~ c(L7, L7)*x7 + c(L8, L8)*x8 + c(L9.group1, L9.group2)*x9 '
## user-defined parameter:
DIF_L9 := L9.group1 - L9.group2
```

Note that this hypothesis is easily tested without a user-defined parameter by using the `lavTestWald()` function. However, a user-defined parameter additionally provides an estimate of the parameter being tested.

User-defined parameters are particularly useful for specifying indirect effects in models of mediation. For example:

```
model <- ' # direct effect
          Y ~ c*X
          # mediator
          M ~ a*X
          Y ~ b*M

          # user defined parameters:

          # indirect effect (a*b)
          ab := a*b
          # total effect (defined using another user-defined parameter)
          total := ab + c
          ,
```

References

Rosseel, Y. (2012). lavaan: An R package for structural equation modeling. *Journal of Statistical Software*, 48(2), 1–36. doi:10.18637/jss.v048.i02

modificationIndices *Modification Indices*

Description

Given a fitted lavaan object, compute the modification indices (= univariate score tests) for a selected set of fixed-to-zero parameters.

Usage

```
modificationIndices(object, standardized = TRUE, cov.std = TRUE,
  information = "expected",
  power = FALSE, delta = 0.1, alpha = 0.05,
  high.power = 0.75, sort. = FALSE, minimum.value = 0,
  maximum.number = nrow(LIST), free.remove = TRUE,
  na.remove = TRUE, op = NULL)
modindices(object, standardized = TRUE, cov.std = TRUE, information = "expected",
  power = FALSE, delta = 0.1, alpha = 0.05, high.power = 0.75,
  sort. = FALSE, minimum.value = 0,
  maximum.number = nrow(LIST), free.remove = TRUE,
  na.remove = TRUE, op = NULL)
```

Arguments

object	An object of class lavaan .
standardized	If TRUE, two extra columns (sepc.lv and sepc.all) will contain standardized values for the EPCs. In the first column (sepc.lv), standardization is based on the variances of the (continuous) latent variables. In the second column (sepc.all), standardization is based on both the variances of both (continuous) observed and latent variables. (Residual) covariances are standardized using (residual) variances.
cov.std	Logical. See standardizedSolution .
information	character indicating the type of information matrix to use (check lavInspect for available options). "expected" information is the default, which provides better control of Type I errors.
power	If TRUE, the (post-hoc) power is computed for each modification index, using the values of delta and alpha.
delta	The value of the effect size, as used in the post-hoc power computation, currently using the unstandardized metric of the epc column.
alpha	The significance level used for deciding if the modification index is statistically significant or not.

high.power	If the computed power is higher than this cutoff value, the power is considered 'high'. If not, the power is considered 'low'. This affects the values in the 'decision' column in the output.
sort.	Logical. If TRUE, sort the output using the values of the modification index values. Higher values appear first.
minimum.value	Numeric. Filter output and only show rows with a modification index value equal or higher than this minimum value.
maximum.number	Integer. Filter output and only show the first maximum number rows. Most useful when combined with the sort. option.
free.remove	Logical. If TRUE, filter output by removing all rows corresponding to free (unconstrained) parameters in the original model.
na.remove	Logical. If TRUE, filter output by removing all rows with NA values for the modification indices.
op	Character string. Filter the output by selecting only those rows with operator op.

Details

Modification indices are just 1-df (or univariate) score tests. The modification index (or score test) for a single parameter reflects (approximately) the improvement in model fit (in terms of the chi-square test statistic), if we would refit the model but allow this parameter to be free. This function is a convenience function in the sense that it produces a (hopefully sensible) table of currently fixed-to-zero (or fixed to another constant) parameters. For each of these parameters, a modification index is computed, together with an expected parameter change (epc) value. It is important to realize that this function will only consider fixed-to-zero parameters. If you have equality constraints in the model, and you wish to examine what happens if you release all (or some) of these equality constraints, use the [lavTestScore](#) function.

Value

A data.frame containing modification indices and EPC's.

Examples

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

fit <- cfa(HS.model, data=HolzingerSwineford1939)
modindices(fit, minimum.value = 10, sort = TRUE)
```

mplus2lavaan

mplus to lavaan converter

Description

Read in an Mplus input file, convert it to lavaan syntax, and fit the model.

Usage

```
mplus2lavaan(inpfile, run = TRUE)
```

Arguments

inpfile	The filename (including a full path) of the Mplus input file. The data (as referred to in the Mplus input file) should be in the same directory as the Mplus input file.
run	Whether to run the specified Mplus input syntax (TRUE) or only to parse and convert the syntax (FALSE).

Value

A lavaan object with the fitted results of the Mplus model. The parsed and converted Mplus syntax is preserved in the @external slot of the lavaan object in the \$mplus.inp element. If run is FALSE, a list of converted syntax is returned.

Author(s)

Michael Hallquist

See Also

[lavExport](#)

Examples

```
## Not run:  
out <- mplus2lavaan("ex5.1.inp")  
summary(out)  
  
## End(Not run)
```

mplus2lavaan.modelSyntax

Convert Mplus model syntax to lavaan

Description

Converts Mplus model syntax into lavaan model syntax.

Usage

```
mplus2lavaan.modelSyntax(syntax)
```

Arguments

`syntax` A character vector containing Mplus model syntax to be converted to lavaan model syntax. Note that parsing Mplus syntax often requires correct usage of newline characters. If `syntax` is a vector of multiple strings, these will be joined with newlines prior to conversion. Alternatively, `\n` characters can be included inline in `syntax`.

Value

A character string of converted lavaan model syntax.

Author(s)

Michael Hallquist

See Also

[mplus2lavaan](#)

Examples

```
## Not run:
syntax <- '
  f1 BY x1*1 x2 x3;
  x1 WITH x2;
  x3 (1);
  x2 (1);
'

lavSyntax <- mplus2lavaan.modelSyntax(syntax)
cat(lavSyntax)

## End(Not run)
```

parameterEstimates *Parameter Estimates*

Description

Parameter estimates of a latent variable model.

Usage

```
parameterEstimates(object,
  se = TRUE, zstat = TRUE, pvalue = TRUE, ci = TRUE,
  standardized = FALSE,
  fmi = FALSE, level = 0.95, boot.ci.type = "perc",
  cov.std = TRUE, fmi.options = list(),
  rsquare = FALSE,
```

```
remove.system.eq = TRUE, remove.eq = TRUE,
remove.ineq = TRUE, remove.def = FALSE,
remove.nonfree = FALSE, remove.step1 = TRUE,
remove.unused = FALSE, add.attributes = FALSE,
output = "data.frame", header = FALSE)
```

Arguments

object	An object of class <code>lavaan</code> .
se	Logical. If TRUE, include column containing the standard errors. If FALSE, this implies <code>zstat</code> and <code>pvalue</code> and <code>ci</code> are also FALSE.
zstat	Logical. If TRUE, an extra column is added containing the so-called z-statistic, which is simply the value of the estimate divided by its standard error.
pvalue	Logical. If TRUE, an extra column is added containing the pvalues corresponding to the z-statistic, evaluated under a standard normal distribution.
ci	If TRUE, confidence intervals are added to the output
level	The confidence level required.
boot.ci.type	If bootstrapping was used, the type of interval required. The value should be one of "norm", "basic", "perc", or "bca.simple". For the first three options, see the help page of the <code>boot.ci</code> function in the <code>boot</code> package. The "bca.simple" option produces intervals using the adjusted bootstrap percentile (BCa) method, but with no correction for acceleration (only for bias). Note that the p-value is still computed assuming that the z-statistic follows a standard normal distribution.
standardized	Logical or character. If TRUE, standardized estimates are added to the output. Note that <i>SEs</i> and tests are still based on unstandardized estimates. Use <code>standardizedSolution</code> to obtain <i>SEs</i> and test statistics for standardized estimates. If a character vector is passed with any of <code>c("std.lv", "std.all", "std.noX")</code> , only the selected standardization methods are added.
cov.std	Logical. If TRUE, the (residual) observed covariances are scaled by the square root of the 'Theta' diagonal elements, and the (residual) latent covariances are scaled by the square root of the 'Psi' diagonal elements. If FALSE, the (residual) observed covariances are scaled by the square root of the diagonal elements of the observed model-implied covariance matrix (Sigma), and the (residual) latent covariances are scaled by the square root of diagonal elements of the model-implied covariance matrix of the latent variables.
fmi	Logical. If TRUE, an extra column is added containing the fraction of missing information for each estimated parameter. Only available if <code>estimator="ML"</code> , <code>missing="(fi)ml"</code> , and <code>se="standard"</code> . See references for more information.
fmi.options	List. If non-empty, arguments can be provided to alter the default options when the model is fitted with the complete(d) data; otherwise, the same options are used as the original model.
remove.eq	Logical. If TRUE, filter the output by removing all rows containing user-specified equality constraints, if any.

<code>remove.system.eq</code>	Logical. If TRUE, filter the output by removing all rows containing system-generated equality constraints, if any.
<code>remove.ineq</code>	Logical. If TRUE, filter the output by removing all rows containing inequality constraints, if any.
<code>remove.def</code>	Logical. If TRUE, filter the output by removing all rows containing parameter definitions, if any.
<code>remove.nonfree</code>	Logical. If TRUE, filter the output by removing all rows containing fixed (non-free) parameters.
<code>remove.step1</code>	Logical. Only used by <code>sam()</code> . If TRUE, filter the output by removing all rows corresponding to the measurement parameters that are part of the first step.
<code>remove.unused</code>	Logical. If TRUE, filter the output by removing all rows containing automatically added parameters (<code>user == 0</code>) that are nonfree, and with their final (<code>est</code>) values fixed to their default values (typically 1 or 0); currently only used for intercepts and scaling-factors.
<code>rsquare</code>	Logical. If TRUE, add additional rows containing the <code>rsquare</code> values (in the <code>est</code> column) of all endogenous variables in the model. Both the <code>lhs</code> and <code>rhs</code> column contain the name of the endogenous variable, while the <code>op</code> column contains <code>r2</code> , to indicate that the values in the <code>est</code> column are <code>rsquare</code> values.
<code>add.attributes</code>	Deprecated argument. Please use <code>output=</code> instead.
<code>output</code>	Character. If <code>"data.frame"</code> , the parameter table is displayed as a standard (albeit lavaan-formatted) <code>data.frame</code> . If <code>"text"</code> (or alias <code>"pretty"</code>), the parameter table is prettyfied, and displayed with subsections (as used by the <code>summary</code> function).
<code>header</code>	Logical. Only used if <code>output = "text"</code> . If TRUE, print a header at the top of the parameter list. This header contains information about the information matrix, if saturated (h1) model is structured or unstructured, and which type of standard errors are shown in the output.

Value

A `data.frame` containing the estimated parameters, parameters, standard errors, and (by default) z-values, p-values, and the lower and upper values of the confidence intervals. If requested, extra columns are added with standardized versions of the parameter estimates.

References

Savalei, V. & Rhemtulla, M. (2012). On obtaining estimates of the fraction of missing information from FIML. *Structural Equation Modeling: A Multidisciplinary Journal*, 19(3), 477-494.

Examples

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

fit <- cfa(HS.model, data=HolzingerSwineford1939)
```

```
parameterEstimates(fit)
parameterEstimates(fit, output = "text")
```

parTable

Parameter Table

Description

Show the parameter table of a fitted model.

Usage

```
parameterTable(object)
parTable(object)
```

Arguments

object An object of class [lavaan](#).

Value

A data.frame containing the model parameters. This is simply the output of the [lavaanify](#) function coerced to a data.frame (with `stringsAsFactors = FALSE`).

See Also

[lavaanify](#).

Examples

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

```
fit <- cfa(HS.model, data=HolzingerSwineford1939)
parTable(fit)
```

PoliticalDemocracy *Industrialization And Political Democracy Dataset*

Description

The ‘famous’ Industrialization and Political Democracy dataset. This dataset is used throughout Bollen’s 1989 book (see pages 12, 17, 36 in chapter 2, pages 228 and following in chapter 7, pages 321 and following in chapter 8). The dataset contains various measures of political democracy and industrialization in developing countries.

Usage

```
data(PoliticalDemocracy)
```

Format

A data frame of 75 observations of 11 variables.

- y1 Expert ratings of the freedom of the press in 1960
- y2 The freedom of political opposition in 1960
- y3 The fairness of elections in 1960
- y4 The effectiveness of the elected legislature in 1960
- y5 Expert ratings of the freedom of the press in 1965
- y6 The freedom of political opposition in 1965
- y7 The fairness of elections in 1965
- y8 The effectiveness of the elected legislature in 1965
- x1 The gross national product (GNP) per capita in 1960
- x2 The inanimate energy consumption per capita in 1960
- x3 The percentage of the labor force in industry in 1960

Source

The dataset was originally retrieved from <http://web.missouri.edu/~kolenikovs/Stat9370/democindus.txt> (link no longer valid; see discussion on SEMNET 18 Jun 2009). The dataset is part of a larger (public) dataset (ICPSR 2532), see <https://www.icpsr.umich.edu/web/ICPSR/studies/2532>.

References

- Bollen, K. A. (1989). *Structural Equations with Latent Variables*. Wiley Series in Probability and Mathematical Statistics. New York: Wiley.
- Bollen, K. A. (1979). Political democracy and the timing of development. *American Sociological Review*, 44, 572-587.
- Bollen, K. A. (1980). Issues in the comparative measurement of political democracy. *American Sociological Review*, 45, 370-390.

Examples

```
head(PoliticalDemocracy)
```

sam

*Fit Structural Equation Models using the SAM approach***Description**

Fit a Structural Equation Model (SEM) using the Structural After Measurement (SAM) approach.

Usage

```

sam(model = NULL, data = NULL, cmd = "sem", se = "twostep",
     mm.list = NULL, mm.args = list(bounds = "wide.zerovar"),
     struc.args = list(estimator = "ML"),
     sam.method = "local", ...,
     local.options = list(M.method = "ML", lambda.correction = TRUE,
                          alpha.correction = 0L, twolevel.method = "h1"),
     global.options = list(), output = "lavaan")

```

Arguments

model	A description of the user-specified model. Typically, the model is described using the lavaan model syntax. See model.syntax for more information. Alternatively, a parameter table (eg. the output of the lavaanify() function) is also accepted.
data	A data frame containing the observed variables used in the model.
cmd	Character. Which command is used to run the sem models. The possible choices are "sem", "cfa" or "lavaan", determining how we deal with default options.
se	Character. The type of standard errors that are used in the final (structural) model. If "twostep" (the default), the standard errors take the estimation uncertainty of the first (measurement) stage into account. If "standard", this uncertainty is ignored, and we treat the measurement information as known. If "none", no standard errors are computed.
mm.list	List. Define the measurement blocks. Each element of the list should be either a single name of a latent variable, or a vector of latent variable names. If omitted, a separate measurement block is used for each latent variable.
mm.args	List. Optional arguments for the fitting function(s) of the measurement block(s) only. See lavOptions for a complete list.
struc.args	List. Optional arguments for the fitting function of the structural part only. See lavOptions for a complete list.
sam.method	Character. Can be set to "local", "global" or "fsr". In the latter case, the results are the same as if Bartlett factor scores were used, without any bias correction.

...	Many more additional options can be defined, using 'name = value'. See lavOptions for a complete list. These options affect both the measurement blocks and the structural part.
local.options	List. Options specific for local SAM method (these options may change over time). If <code>lambda.correction = TRUE</code> , we ensure that the variance matrix of the latent variables (VETA) is positive definite. The <code>alpha.correction</code> options must be an integer. Acceptable values are in the range 0 till N-1. If zero (the default), no small sample correction is performed, and the bias-correction is the same as with local SAM. When equal to N-1, the bias-correction is eliminated, and the results are the same as naive FSR. Typical values are 0, P+1 (where P is the number of predictors in the structural model), P+5, and (N-1)/2.
global.options	List. Options specific for global SAM method (not used for now).
output	Character. If "lavaan", a lavaan object returned. If "list", a list is returned with all the ingredients from the different stages.

Details

The `sam` function tries to automate the SAM approach, by first estimating the measurement part of the model, and then the structural part of the model. See reference for more details.

Note that in the current implementation, all indicators of latent variables have to be observed. This implies: no support for second-order factor structures (for now).

Value

If `output = "lavaan"`, an object of class `lavaan`, for which several methods are available, including a summary method. If `output = "list"`, a list.

References

Rosseel and Loh (2021). A structural-after-measurement approach to Structural Equation Modeling. *Psychological Methods*. Advance online publication. <https://dx.doi.org/10.1037/met0000503>

See Also

[lavaan](#)

Examples

```
## The industrialization and Political Democracy Example
## Bollen (1989), page 332
model <- '
  # latent variable definitions
  ind60 =~ x1 + x2 + x3
  dem60 =~ y1 + a*y2 + b*y3 + c*y4
  dem65 =~ y5 + a*y6 + b*y7 + c*y8

  # regressions
  dem60 ~ ind60
  dem65 ~ ind60 + dem60
```



```

# residual correlations
y1 ~~ y5
y2 ~~ y4 + y6
y3 ~~ y7
y4 ~~ y8
y6 ~~ y8
,

fit.sam <- sam(model, data = PoliticalDemocracy,
              mm.list = list(ind = "ind60", dem = c("dem60", "dem65")))
summary(fit.sam)

```

sem

*Fit Structural Equation Models***Description**

Fit a Structural Equation Model (SEM).

Usage

```

sem(model = NULL, data = NULL, ordered = NULL, sampling.weights = NULL,
     sample.cov = NULL, sample.mean = NULL, sample.th = NULL,
     sample.nobs = NULL, group = NULL, cluster = NULL,
     constraints = "", WLS.V = NULL, NACOV = NULL, ov.order = "model",
     ...)

```

Arguments

model	A description of the user-specified model. Typically, the model is described using the lavaan model syntax. See model.syntax for more information. Alternatively, a parameter table (eg. the output of the <code>lavaanify()</code> function) is also accepted.
data	An optional data frame containing the observed variables used in the model. If some variables are declared as ordered factors, lavaan will treat them as ordinal variables.
ordered	Character vector. Only used if the data is in a data.frame. Treat these variables as ordered (ordinal) variables, if they are endogenous in the model. Importantly, all other variables will be treated as numeric (unless they are declared as ordered in the data.frame.) Since 0.6-4, ordered can also be logical. If TRUE, all observed endogenous variables are treated as ordered (ordinal). If FALSE, all observed endogenous variables are considered to be numeric (again, unless they are declared as ordered in the data.frame.)
sampling.weights	A variable name in the data frame containing sampling weight information. Currently only available for non-clustered data. Depending on the <code>sampling.weights.normalization</code> option, these weights may be rescaled (or not) so that their sum equals the number of observations (total or per group).

<code>sample.cov</code>	Numeric matrix. A sample variance-covariance matrix. The rownames and/or colnames must contain the observed variable names. For a multiple group analysis, a list with a variance-covariance matrix for each group.
<code>sample.mean</code>	A sample mean vector. For a multiple group analysis, a list with a mean vector for each group.
<code>sample.th</code>	Vector of sample-based thresholds. For a multiple group analysis, a list with a vector of thresholds for each group.
<code>sample.nobs</code>	Number of observations if the full data frame is missing and only sample moments are given. For a multiple group analysis, a list or a vector with the number of observations for each group.
<code>group</code>	Character. A variable name in the data frame defining the groups in a multiple group analysis.
<code>cluster</code>	Character. A (single) variable name in the data frame defining the clusters in a two-level dataset.
<code>constraints</code>	Additional (in)equality constraints not yet included in the model syntax. See model.syntax for more information.
<code>WLS.V</code>	A user provided weight matrix to be used by estimator "WLS"; if the estimator is "DWLS", only the diagonal of this matrix will be used. For a multiple group analysis, a list with a weight matrix for each group. The elements of the weight matrix should be in the following order (if all data is continuous): first the means (if a meanstructure is involved), then the lower triangular elements of the covariance matrix including the diagonal, ordered column by column. In the categorical case: first the thresholds (including the means for continuous variables), then the slopes (if any), the variances of continuous variables (if any), and finally the lower triangular elements of the correlation/covariance matrix excluding the diagonal, ordered column by column.
<code>NACOV</code>	A user provided matrix containing the elements of (N times) the asymptotic variance-covariance matrix of the sample statistics. For a multiple group analysis, a list with an asymptotic variance-covariance matrix for each group. See the <code>WLS.V</code> argument for information about the order of the elements.
<code>ov.order</code>	Character. If "model" (the default), the order of the observed variable names (as reflected for example in the output of <code>lavNames()</code>) is determined by the model syntax. If "data", the order is determined by the data (either the full data.frame or the sample (co)variance matrix). If the <code>WLS.V</code> and/or <code>NACOV</code> matrices are provided, this argument is currently set to "data".
<code>...</code>	Many more additional options can be defined, using 'name = value'. See lavOptions for a complete list.

Details

The `sem` function is a wrapper for the more general [lavaan](#) function, but setting the following default options: `int.ov.free = TRUE`, `int.lv.free = FALSE`, `auto.fix.first = TRUE` (unless `std.lv = TRUE`), `auto.fix.single = TRUE`, `auto.var = TRUE`, `auto.cov.lv.x = TRUE`, `auto.efa = TRUE`, `auto.th = TRUE`, `auto.delta = TRUE`, and `auto.cov.y = TRUE`.

Value

An object of class `lavaan`, for which several methods are available, including a summary method.

References

Yves Rosseel (2012). `lavaan`: An R Package for Structural Equation Modeling. *Journal of Statistical Software*, 48(2), 1-36. doi:[10.18637/jss.v048.i02](https://doi.org/10.18637/jss.v048.i02)

See Also

[lavaan](#)

Examples

```
## The industrialization and Political Democracy Example
## Bollen (1989), page 332
model <- '
  # latent variable definitions
  ind60 =~ x1 + x2 + x3
  dem60 =~ y1 + a*y2 + b*y3 + c*y4
  dem65 =~ y5 + a*y6 + b*y7 + c*y8

  # regressions
  dem60 ~ ind60
  dem65 ~ ind60 + dem60

  # residual correlations
  y1 ~~ y5
  y2 ~~ y4 + y6
  y3 ~~ y7
  y4 ~~ y8
  y6 ~~ y8
'

fit <- sem(model, data = PoliticalDemocracy)
summary(fit, fit.measures = TRUE)
```

Description

Simulate data starting from a lavaan model syntax.

Usage

```
simulateData(model = NULL, model.type = "sem", meanstructure = FALSE,
  int.ov.free = TRUE, int.lv.free = FALSE,
  marker.int.zero = FALSE, conditional.x = FALSE, fixed.x = FALSE,
  orthogonal = FALSE, std.lv = TRUE, auto.fix.first = FALSE,
  auto.fix.single = FALSE, auto.var = TRUE, auto.cov.lv.x = TRUE,
  auto.cov.y = TRUE, ..., sample.nobs = 500L, ov.var = NULL,
  group.label = paste("G", 1:ngroups, sep = ""), skewness = NULL,
  kurtosis = NULL, seed = NULL, empirical = FALSE,
  return.type = "data.frame", return.fit = FALSE,
  debug = FALSE, standardized = FALSE)
```

Arguments

model	A description of the user-specified model. Typically, the model is described using the lavaan model syntax. See model.syntax for more information. Alternatively, a parameter table (eg. the output of the lavaanify() function) is also accepted.
model.type	Set the model type: possible values are "cfa", "sem" or "growth". This may affect how starting values are computed, and may be used to alter the terminology used in the summary output, or the layout of path diagrams that are based on a fitted lavaan object.
meanstructure	If TRUE, the means of the observed variables enter the model. If "default", the value is set based on the user-specified model, and/or the values of other arguments.
int.ov.free	If FALSE, the intercepts of the observed variables are fixed to zero.
int.lv.free	If FALSE, the intercepts of the latent variables are fixed to zero.
marker.int.zero	Logical. Only relevant if the metric of each latent variable is set by fixing the first factor loading to unity. If TRUE, it implies meanstructure = TRUE and std.lv = FALSE, and it fixes the intercepts of the marker indicators to zero, while freeing the means/intercepts of the latent variables. Only works correctly for single group, single level models.
conditional.x	If TRUE, we set up the model conditional on the exogenous 'x' covariates; the model-implied sample statistics only include the non-x variables. If FALSE, the exogenous 'x' variables are modeled jointly with the other variables, and the model-implied statistics reflect both sets of variables. If "default", the value is set depending on the estimator, and whether or not the model involves categorical endogenous variables.
fixed.x	If TRUE, the exogenous 'x' covariates are considered fixed variables and the means, variances and covariances of these variables are fixed to their sample values. If FALSE, they are considered random, and the means, variances and covariances are free parameters. If "default", the value is set depending on the mimic option.
orthogonal	If TRUE, the exogenous latent variables are assumed to be uncorrelated.

std.lv	If TRUE, the metric of each latent variable is determined by fixing their variances to 1.0. If FALSE, the metric of each latent variable is determined by fixing the factor loading of the first indicator to 1.0.
auto.fix.first	If TRUE, the factor loading of the first indicator is set to 1.0 for every latent variable.
auto.fix.single	If TRUE, the residual variance (if included) of an observed indicator is set to zero if it is the only indicator of a latent variable.
auto.var	If TRUE, the (residual) variances of both observed and latent variables are set free.
auto.cov.lv.x	If TRUE, the covariances of exogenous latent variables are included in the model and set free.
auto.cov.y	If TRUE, the covariances of dependent variables (both observed and latent) are included in the model and set free.
...	additional arguments passed to the lavaan function.
sample.nobs	Number of observations. If a vector, multiple datasets are created. If <code>return.type = "matrix"</code> or <code>return.type = "cov"</code> , a list of <code>length(sample.nobs)</code> is returned, with either the data or covariance matrices, each one based on the number of observations as specified in <code>sample.nobs</code> . If <code>return.type = "data.frame"</code> , all datasets are merged and a group variable is added to mimic a multiple group dataset.
ov.var	The user-specified variances of the observed variables.
group.label	The group labels that should be used if multiple groups are created.
skewness	Numeric vector. The skewness values for the observed variables. Defaults to zero.
kurtosis	Numeric vector. The kurtosis values for the observed variables. Defaults to zero.
seed	Set random seed.
empirical	Logical. If TRUE, the implied moments (Mu and Sigma) specify the empirical not population mean and covariance matrix.
return.type	If <code>"data.frame"</code> , a <code>data.frame</code> is returned. If <code>"matrix"</code> , a numeric matrix is returned (without any variable names). If <code>"cov"</code> , a covariance matrix is returned (without any variable names).
return.fit	If TRUE, return the fitted model that has been used to generate the data as an attribute (called <code>"fit"</code>); this may be useful for inspection.
debug	If TRUE, debugging information is displayed.
standardized	If TRUE, the residual variances of the observed variables are set in such a way such that the model implied variances are unity. This allows regression coefficients and factor loadings (involving observed variables) to be specified in a standardized metric.

Details

Model parameters can be specified by fixed values in the lavaan model syntax. If no fixed values are specified, the value zero will be assumed, except for factor loadings and variances, which are

set to unity by default. By default, multivariate normal data are generated. However, by providing skewness and/or kurtosis values, nonnormal multivariate data can be generated, using the Vale & Maurelli (1983) method.

Value

The generated data. Either as a data.frame (if `return.type="data.frame"`), a numeric matrix (if `return.type="matrix"`), or a covariance matrix (if `return.type="cov"`).

Examples

```
# specify population model
population.model <- ' f1 =~ x1 + 0.8*x2 + 1.2*x3
                    f2 =~ x4 + 0.5*x5 + 1.5*x6
                    f3 =~ x7 + 0.1*x8 + 0.9*x9

                    f3 ~ 0.5*f1 + 0.6*f2
                    '

# generate data
set.seed(1234)
myData <- simulateData(population.model, sample.nobs=100L)

# population moments
fitted(sem(population.model))

# sample moments
round(cov(myData), 3)
round(colMeans(myData), 3)

# fit model
myModel <- ' f1 =~ x1 + x2 + x3
            f2 =~ x4 + x5 + x6
            f3 =~ x7 + x8 + x9
            f3 ~ f1 + f2 '
fit <- sem(myModel, data=myData)
summary(fit)
```

standardizedSolution *Standardized Solution*

Description

Standardized solution of a latent variable model.

Usage

```
standardizedSolution(object, type = "std.all", se = TRUE, zstat = TRUE,
                    pvalue = TRUE, ci = TRUE, level = 0.95, cov.std = TRUE,
                    remove.eq = TRUE, remove.ineq = TRUE, remove.def = FALSE,
```

```
partable = NULL, GLIST = NULL, est = NULL,
output = "data.frame")
```

Arguments

object	An object of class <code>lavaan</code> .
type	If <code>"std.lv"</code> , the standardized estimates are on the variances of the (continuous) latent variables only. If <code>"std.all"</code> , the standardized estimates are based on both the variances of both (continuous) observed and latent variables. If <code>"std.nox"</code> , the standardized estimates are based on both the variances of both (continuous) observed and latent variables, but not the variances of exogenous covariates.
se	Logical. If TRUE, standard errors for the standardized parameters will be computed, together with a z-statistic and a p-value.
zstat	Logical. If TRUE, an extra column is added containing the so-called z-statistic, which is simply the value of the estimate divided by its standard error.
pvalue	Logical. If TRUE, an extra column is added containing the pvalues corresponding to the z-statistic, evaluated under a standard normal distribution.
ci	If TRUE, simple symmetric confidence intervals are added to the output
level	The confidence level required.
cov.std	Logical. If TRUE, the (residual) observed covariances are scaled by the square root of the 'Theta' diagonal elements, and the (residual) latent covariances are scaled by the square root of the 'Psi' diagonal elements. If FALSE, the (residual) observed covariances are scaled by the square root of the diagonal elements of the observed model-implied covariance matrix (Sigma), and the (residual) latent covariances are scaled by the square root of diagonal elements of the model-implied covariance matrix of the latent variables.
remove.eq	Logical. If TRUE, filter the output by removing all rows containing equality constraints, if any.
remove.ineq	Logical. If TRUE, filter the output by removing all rows containing inequality constraints, if any.
remove.def	Logical. If TRUE, filter the output by removing all rows containing parameter definitions, if any.
GLIST	List of model matrices. If provided, they will be used instead of the GLIST inside the object@Model slot. Only works if the est argument is also provided. See Note.
est	Numeric. Parameter values (as in the 'est' column of a parameter table). If provided, they will be used instead of the parameters that can be extract from object. Only works if the GLIST argument is also provided. See Note.
partable	A custom list or data.frame in which to store the standardized parameter values. If provided, it will be used instead of the parameter table inside the object@ParTable slot.
output	Character. If <code>"data.frame"</code> , the parameter table is displayed as a standard (albeit lavaan-formatted) data.frame. If <code>"text"</code> (or alias <code>"pretty"</code>), the parameter table is prettyfied, and displayed with subsections (as used by the summary function).

Value

A data.frame containing standardized model parameters.

Note

The est, GLIST, and partable arguments are not meant for everyday users, but for authors of external R packages that depend on lavaan. Only to be used with great caution.

Examples

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

fit <- cfa(HS.model, data=HolzingerSwineford1939)
standardizedSolution(fit)
```

summary.efaList	<i>Summarizing EFA Fits</i>
-----------------	-----------------------------

Description

S3 summary and print methods for class efaList.

Usage

```
## S3 method for class 'efaList'
summary(object,
        nd = 3L, cutoff = 0.3, dot.cutoff = 0.1, alpha.level = 0.01,
        lambda = TRUE, theta = TRUE, psi = TRUE, fit.table = TRUE,
        fs.determinacy = FALSE, eigenvalues = TRUE, sumsq.table = TRUE,
        lambda.structure = FALSE, se = FALSE, zstat = FALSE,
        pvalue = FALSE, ...)

## S3 method for class 'efaList.summary'
print(x, nd = 3L, cutoff = 0.3, dot.cutoff = 0.1,
      alpha.level = 0.01, ...)
```

Arguments

object	An object of class efaList, usually, a result of a call to <code>efa</code> with (the default) <code>output = "efa"</code> .
x	An object of class <code>summary.efaList</code> , usually, a result of a call to <code>summary.efaList</code> .
nd	Integer. The number of digits that are printed after the decimal point in the output.

cutoff	Numeric. Factor loadings smaller than this value (in absolute value) are not printed (even if they are significantly different from zero). The idea is that only medium to large factor loadings are printed, to better see the overall structure.
dot.cutoff	Numeric. Factor loadings larger (in absolute value) than this value, but smaller (in absolute value) than the cutoff value are shown as a dot. They represent small loadings that may still need your attention.
alpha.level	Numeric. If the p-value of a factor loading is smaller than this value, a significance star is printed to the right of the factor loading. To switch this off, use <code>alpha.level = 0</code> .
lambda	Logical. If TRUE, include the (standardized) factor loadings in the summary.
theta	Logical. If TRUE, include the unique variances and the communalities in the table of factor loadings.
psi	Logical. If TRUE, include the factor correlations in the summary. Ignored if only a single factor is used.
fit.table	Logical. If TRUE, show fit information for each model.
fs.determinacy	Logical. If TRUE, show the factor score determinacy values per factor (assuming regression factor scores are used) and their squared values.
eigenvalues	Logical. If TRUE, include the eigenvalues of the sample variance-covariance matrix in the summary.
sumsq.table	Logical. If TRUE, include a table including sums of squares of factor loadings (and related measures) in the summary. The sums of squares are computed as the diagonal elements of Lambda times Psi (where Psi is the matrix of factor correlations.). If orthogonal rotation was used, Psi is diagonal and the sums of squares are identical to the sums of the squared column elements of the Lambda matrix (i.e., the factor loadings). This is no longer the case when oblique rotation has been used. But in both cases (orthogonal or oblique), the (total) sum of the sums of squares equals the sum of the communalities. In the second row of the table (Proportion of total), the sums of squares are divided by the total. In the third row of the table (Proportion var), the sums of squares are divided by the number of items.
lambda.structure	Logical. If TRUE, show the structure matrix (i.e., the factor loadings multiplied by the factor correlations).
se	Logical. If TRUE, include the standard errors of the standardized lambda, theta and psi elements in the summary.
zstat	Logical. If TRUE, include the Z-statistics of the standardized lambda, theta and psi elements in the summary.
pvalue	Logical. If TRUE, include the P-values of the standardized lambda, theta and psi elements in the summary.
...	Further arguments passed to or from other methods.

Value

The function `summary.efaList` computes and returns a list of summary statistics for the list of EFA models in object.

Examples

```
## The famous Holzinger and Swineford (1939) example
fit <- efa(data = HolzingerSwineford1939,
           ov.names = paste("x", 1:9, sep = ""),
           nfactors = 1:3,
           rotation = "geomin",
           rotation.args = list(geomin.epsilon = 0.01, rstarts = 1))
summary(fit, nd = 3L, cutoff = 0.2, dot.cutoff = 0.05,
        lambda.structure = TRUE, pvalue = TRUE)
```

varTable

*Variable Table***Description**

Summary information about the variables included in either a data.frame, or a fitted lavaan object.

Usage

```
varTable(object, ov.names = names(object), ov.names.x = NULL,
         ordered = NULL, factor = NULL, as.data.frame. = TRUE)
```

Arguments

object	Either a data.frame, or an object of class lavaan .
ov.names	Only used if object is a data.frame. A character vector containing the variables that need to be summarized.
ov.names.x	Only used if object is a data.frame. A character vector containing additional variables that need to be summarized.
ordered	Character vector. Which variables should be treated as ordered factors
factor	Character vector. Which variables should be treated as (unordered) factors?
as.data.frame.	If TRUE, return the list as a data.frame.

Value

A list or data.frame containing summary information about variables in a data.frame. If object is a fitted lavaan object, it displays the summary information about the observed variables that are included in the model. The summary information includes variable type (numeric, ordered, ...), the number of non-missing values, the mean and variance for numeric variables, the number of levels of ordered variables, and the labels for ordered variables.

Examples

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

fit <- cfa(HS.model, data=HolzingerSwineford1939)
varTable(fit)
```

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