

Package ‘asremlPlus’

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Title Augments the Use of 'Asreml' in Fitting Mixed Models

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Depends R (>= 2.10.0)

Imports dae, ggplot2, stats, methods, utils, grDevices

Enhances asreml

SystemRequirements asreml-R 3.x

Description Assists in automating the testing of terms in mixed models when 'asreml' is used to fit the models. The content falls into the following natural groupings: (i) Data, (ii) Object manipulation functions, (iii) Model modification functions, (iv) Model testing functions, (v) Model diagnostics functions, (vi) Prediction production and presentation functions, (vii) Response transformation functions, and (viii) Miscellaneous functions. A history of the fitting of a sequence of models is kept in a data frame. Procedures are available for choosing models that conform to the hierarchy or marginality principle and for displaying predictions for significant terms in tables and graphs. The package 'asreml' provides a computationally efficient algorithm for fitting mixed models using Residual Maximum Likelihood. It can be purchased from 'VSNi' (<http://www.vsn.co.uk/>) as 'asreml-R', who will supply a zip file for local installation/updating.

License GPL (>=2)

URL <http://chris.brien.name>

NeedsCompilation no

R topics documented:

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| | |
|--------------------|---|
| asremlPlus-package | <i>Augments the use of asreml-R in fitting mixed models</i> |
|--------------------|---|

Description

Assists in automating the testing of terms in mixed models when 'asreml' is used to fit the models. The content falls into the following natural groupings: (i) Data, (ii) Object manipulation functions, (iii) Model modification functions, (iv) Model testing functions, (v) Model diagnostics functions, (vi) Prediction production and presentation functions, (vii) Response transformation functions, and (viii) Miscellaneous functions. A history of the fitting of a sequence of models is kept in a data frame. Procedures are available for choosing models that conform to the hierarchy or marginality principle and for displaying predictions for significant terms in tables and graphs. The package 'asreml' provides a computationally efficient algorithm for fitting mixed models using Residual Maximum Likelihood. It can be purchased from 'VSNi' (<http://www.vsn.co.uk/>) as 'asreml-R', who will supply a zip file for local installation/updating.

Details

The DESCRIPTION file:

| | |
|----------|--|
| Package: | asremlPlus |
| Version: | 2.0-3 |
| Date: | 2015-11-11 |
| Title: | Augments the Use of 'Asreml' in Fitting Mixed Models |
| Author: | Chris Brien <Chris.Brien@unisa.edu.au>. |

Maintainer: Chris Brien <Chris.Brien@unisa.edu.au>
 Depends: R (>= 2.10.0)
 Imports: dae, ggplot2, stats, methods, utils, grDevices
 Enhances: asreml
 SystemRequirements: asreml-R 3.x
 Description: Assists in automating the testing of terms in mixed models when 'asreml' is used to fit the model.
 License: GPL (>=2)
 URL: <http://chris.brien.name>

Index:

(i) Data

Wheat.dat Data for an experiment to investigate 25 varieties of wheat.

(ii) Object manipulation

alldiffs Forms an object of S3-class 'alldiffs' that stores the predictions for a model fitted using asreml, along with statistics for all pairwise differences.

asrtests Forms an object of S3-class 'asrtests' that stores a fitted asreml object, a pseudo-anova table for the fixed and a history of changes and hypothesis testing used in obtaining the model.

print.alldiffs Prints the values in an 'alldiffs' object in a nice format.

print.asrtests Prints the values in an 'asrtests' object

(iii) Model modification

addrm.terms.asrtests Adds or removes the specified set terms from either the fixed or random model and records the change in a data.frame.

newfit.asreml Refits an asreml model with modified model formula using either a call to 'update.asreml' or a direct call to 'asreml'.

newrcov.asrtests Fits a new rcov formula using asreml.

rmboundary.asrtests Removes any boundary or singular variance components from the fit stored in 'asreml.obj' and records their removal in a data.frame.

setvarianceterms.asreml Allows the setting of constraints and initial values for terms in the 'random' and 'rcov' arguments of an 'asreml' call.

sig.devn.reparam.asrtests This function reparameterizes each random (deviations) term involving 'devn.fac' to a fixed term and ensures that the same term, with 'trend.num' replacing 'devn.fac', is included if any other term with 'trend.num' is included in 'terms'.

(iv) Model testing

| | |
|---------------------------------------|---|
| <code>choose.model.asrtests</code> | Determines the set of significant terms taking into account hierarchy or marginality relations. |
| <code>info.crit.asreml</code> | Computes AIC and BIC for a model. |
| <code>recalc.wald.tab.asrtests</code> | Recalculates the denDF, F.inc and P values for a table of Wald test statistics obtained using 'wald.asreml'. |
| <code>reml.lrt.asreml</code> | Performs REML likelihood ratio test. |
| <code>testranfix.asrtests</code> | Tests for a single fixed or random term in model fitted using 'asreml'. |
| <code>testrcov.asrtests</code> | Fits a new rcov formula using 'asreml' and tests whether the change is significant. |
| <code>testswapan.asrtests</code> | Tests, using a REMLRT, whether the difference between current random model and one in which oldterms are dropped and newterms are added is significant. |

(v) Model diagnostics

| | |
|--------------------------------|---|
| <code>variofaces.asreml</code> | Plot empirical variogram faces, including envelopes, as described by Stefanova, Smith & Cullis (2009) |
|--------------------------------|---|

(vi) Prediction production and presentation

| | |
|-------------------------------------|---|
| <code>pred.present.asreml</code> | This function forms the predictions for each significant term and presents them in tables and/or graphs. |
| <code>predictiondiffs.asreml</code> | Forms all pairwise differences between a set of predictions, their standard errors and p-values for a test of whether the differences are significantly different from zero. |
| <code>predictionplot.asreml</code> | This function plots the predictions for a term, possibly with error bars. |
| <code>predictparallel.asreml</code> | This function forms the predictions and associated statistics for a term, taking into account that a numeric vector and a factor having parallel values may occur in the model. It stores the results in an object of class 'alldiffs' and prints the results. It can be used when there are not parallel values. |

(vii) Response transformation

| | |
|------------------------------|--|
| <code>angular</code> | Applies the angular transformation to proportions. |
| <code>angular.mod</code> | Applies the modified angular transformation to a vector of counts. |
| <code>power.transform</code> | Perform linear and power transformations on a variable whose name is given as a character string in 'var.name'. The transformed variable is stored in the 'data.frame data'. |

(viii) Miscellaneous

| | |
|-------------------------|---|
| <code>num.recode</code> | Recodes the unique values of a vector using the values in a new vector. |
|-------------------------|---|

| | |
|--------------------------|--|
| permute.square | Permutes the rows and columns of a square matrix. |
| permute.to.zero.lowertri | Permutes a square matrix until all the lower triangular elements are zero. |

The functions whose names end in 'asrtests', which are most of the model functions, utilize an asrtests object that stores: (i) the currently fitted model in asreml.obj, (ii) the table of test statistics for the fixed effects in wald.tab, and (iii) a data frame that contains a history of the changes made to the model in test.summary.

Author(s)

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References

Butler, D. G., et al. (2010). *Analysis of Mixed Models for S language environments: ASReml-R reference manual*. Brisbane, DPI Publications.

Stefanova, K. T., Smith, A. B. & Cullis, B. R. (2009) Enhanced diagnostics for the spatial analysis of field trials. *Journal of Agricultural, Biological, and Environmental Statistics*, **14**, 392–410.

See Also

asreml

Examples

```
## Not run:
## Analyse wheat dat using asreml and asremlPlus
## Set up for analysis
library(dae)
library(asreml)
## use ?Wheat.dat for data set details
data(Wheat.dat)

# Fit initial model
current.asr <- asreml(yield ~ Rep + WithinColPairs + Variety,
                    random = ~ Row + Column + units,
                    rcov = ~ ar1(Row):ar1(Column),
                    data=Wheat.dat)
summary(current.asr)

# Load current fit into an asrtests object
current.asrt <- asrtests(current.asr, NULL, NULL)

# Check for and remove any boundary terms
current.asrt <- rmboundary.asrtests(current.asrt)

# Check term for within Column pairs
current.asrt <- testranfix.asrtests("WithinColPairs", current.asrt, drop.fix.ns=TRUE)

# Test nugget term
current.asrt <- testranfix.asrtests("units", current.asrt, positive=TRUE)

# Test Row autocorrelation
```

```

current.asrt <- testrcov.asrtests("~ Row:ar1(Column)", current.asrt,
                                label="Row autocorrelation", simpler=TRUE)

# Test Col autocorrelation (depends on whether Row autocorrelation retained)
k <- match("Row autocorrelation", current.asrt$test.summary$terms)
p <- current.asrt$test.summary$p
if (p[k] <= 0.05)
  current.asrt <- testrcov.asrtests("~ ar1(Row):Column", current.asrt,
                                    label="Col autocorrelation", simpler=TRUE,
                                    update=FALSE)
else
  current.asrt <- testrcov.asrtests("~ Row:Column", current.asrt,
                                    label="Col autocorrelation", simpler=TRUE,
                                    update=FALSE)
}
print(current.asrt)

# Get current fitted asreml object
current.asr <- current.asrt$asreml.obj
current.asr <- update(current.asr, aom=TRUE)

# Do residuals-versus-fitted values plot
plot(fitted.values(current.asr),residuals(current.asr))

# Form variance matrix based on estimated variance parameters
s2 <- current.asr$sigma2
gamma.Row <- current.asr$gammas[1]
gamma.unit <- current.asr$gammas[2]
rho.r <- current.asr$gammas[4]
rho.c <- current.asr$gammas[5]
row.ar1 <- mat.ar1(order=10, rho=rho.r)
col.ar1 <- mat.ar1(order=15, rho=rho.c)
V <- fac.vcmat(Wheat.dat$Row, gamma.Row) +
  gamma.unit * diag(1, nrow=150, ncol=150) +
  mat.dirprod(row.ar1, col.ar1)
V <- s2*V

#Produce variogram and variogram faces plot (Stefanaova et al, 2009)
plot.asrVariogram(variogram(current.asr))
variofaces.asreml(current.asr, V=V)

#Get Variety predictionsand all pairwise prediction differences and p-values
Var.diffs <- predictparallel.asreml(classify = "Variety",
                                   asreml.obj=current.asr,
                                   error.intervals="halfLeast",
                                   wald.tab=current.asrt$wald.tab,
                                   tables = "predictions")
print(Var.diffs, which = c("differences", "p.differences"))

## End(Not run)

```

`addrm.terms.asrtests` *Adds or removes the specified set terms from either the fixed or random model and records the change in a data.frame.*

Description

The specified terms are simply added or removed from either the fixed or random model. No hypothesis testing is performed and no check is made for boundary or singular terms. A row is added to the `test.summary.data.frame` stating whether fixed or random terms have been added or removed. Convergence in fitting the model is checked and a note included in the action if there was not. All components of the `asrtests` object are updated, although `wald.tab` is only updated if it is present in the supplied `asrtests` object.

Usage

```
addrm.terms.asrtests(terms = NULL, asrtests.obj, add = FALSE,
                     random = FALSE, label = NULL,
                     denDF = "default", trace = FALSE,
                     update = TRUE, set.terms = NULL, ignore.suffices = TRUE,
                     constraints = "P", initial.values = NA, ...)
```

Arguments

| | |
|------------------------------|--|
| <code>terms</code> | a single character string in the form of a formula which, after expansion, specifies the sum of a set of terms to be added or dropped. |
| <code>asrtests.obj</code> | an <code>asrtests</code> object for a fitted model that is a list containing an <code>asreml</code> object, a <code>wald.tab</code> data.frame with 4 columns, and a data.frame with 5 columns that records any previous changes and tests in the fitted model. |
| <code>add</code> | whether to add or remove terms from the model. |
| <code>random</code> | whether terms are to added or removed from the fixed or random model. |
| <code>label</code> | a character string to use as the label in <code>test.summary</code> and which indicates what is being tested. If <code>label</code> is <code>NULL</code> , either <code>Fixed terms</code> or <code>Random terms</code> is used, depending on whether <code>random</code> is <code>TRUE</code> or <code>FALSE</code> . |
| <code>denDF</code> | Specifies the enthod to use in computing approximate denominator degrees of freedom when <code>wald.asreml</code> is called. Can be <code>none</code> to suppress the computations, <code>numeric</code> for numerical methods, <code>algebraic</code> for algebraic methods or <code>default</code> , the default, to autommatically choose numeric or algebraic computations depending on problem size. The denominator degrees of freedom are calculated according to Kenward and Roger (1997) for fixed terms in the dense part of the model. |
| <code>trace</code> | if <code>TRUE</code> then partial iteration details are displayed when <code>ASReml-R</code> functions are invoked; if <code>FALSE</code> then no output is displayed. |
| <code>update</code> | if <code>TRUE</code> then <code>update.asreml</code> is called in removing and adding terms to the model. In doing this the arguments <code>R.param</code> and <code>G.param</code> are set to those in the <code>asreml</code> object stored in the supplied <code>asrtests.obj</code> so that the values from the previous model are used as starting values. If <code>FALSE</code> then calls are made to <code>asreml</code> in which the only changes from the previous call are (i) that the random model is updated and (ii) modifications specified via <code>...</code> are made. |
| <code>set.terms</code> | a character vector specifying the terms that are to have constraints and/or initial values set prior to fitting. |
| <code>ignore.suffices</code> | a logical vector specifying whether the suffices of the <code>asreml</code> -assigned names of the variance terms (i.e. the information to the right of an "!", other than "R!") is to be ignored in matching elements of terms. If <code>TRUE</code> for an element of terms, the suffices are stripped from the <code>asreml</code> -assigned names. If <code>FALSE</code> |

| | |
|-----------------------------|---|
| | for an element of <code>terms</code> , the element must exactly match an <code>asreml</code> -assigned name for a variance term. This vector must be of length one or the same length as <code>terms</code> . If it is of length one then the same action is applied to the <code>asreml</code> -assigned suffices for all the terms in <code>terms</code> . |
| <code>constraints</code> | a character vector specifying the constraints to be applied to the terms specified in <code>terms</code> . This vector must be of length one or the same length as <code>terms</code> . If it is of length one then the same constraint is applied to all the terms in <code>terms</code> . If any of the constraints are equal to <code>NA</code> then they are left unchanged for those terms. |
| <code>initial.values</code> | a character vector specifying the initial values for the terms specified in <code>terms</code> . This vector must be of length one or the same length as <code>terms</code> . If it is of length one then the same initial value is applied to all the terms in <code>terms</code> . If any of the <code>initial.values</code> are equal to <code>NA</code> then they are left unchanged for those terms. |
| <code>...</code> | further arguments passed to <code>asreml</code> and to <code>wald.asreml</code> . |

Value

An `asrtests` object, which is a list containing:

1. `asreml.obj`: an `asreml` object containing the fit of the model after all boundary and singular terms have been removed;
2. `wald.tab`: a 4-column data.frame containing a pseudo-anova table for the fixed terms produced by `wald.asreml`;
3. `test.summary`: a data.frame with columns `term`, `DF`, `denDF`, `p` and `action`. A row is added to it for each term that is dropped, added or tested or a note that several terms have been added or removed. A row contains the name of the term, the DF, the p-value and the action taken. Possible codes are: Dropped, Retained, Swapped, Unswapped, Significant, Nonsignificant, Absent, Added, Removed and Boundary. If the changed model did not converge, Unconverged will be added to the code. Note that the logical `asreml.obj$converge` also reflects whether there is convergence.

See Also

[asrtests](#), [rmboundary.asrtests](#), [testranfix.asrtests](#), [testrcov.asrtests](#), [newfit.asreml](#), [sig.devn.reparam.asrtests](#), [choose.model.asrtests](#)

Examples

```
## Not run:
terms <- "(Date/(Sources * (Type + Species)))"
current.asrt <- addrm.terms.asrtests(terms, current.asrt, add = TRUE)

current.asrt <- addrm.terms.asrtests("A + B", current.asrt, denDF = "algebraic")

## End(Not run)
```

| | |
|----------|---|
| alldiffs | <i>Forms an object of S3-class alldiffs that stores the predictions for a model fitted using asreml, along with supplied statistics for all pairwise differences.</i> |
|----------|---|

Description

Creates an object of S3-class alldiffs that consists of a list containing the following components: predictions, differences, p.differences, sed, LSD and backtransforms. Predictions must be supplied to the functions while the others will be set only if they are supplies; those not supplied are set to NULL. It also has attributes response, response.title, term, classify and tdf, which will be set to the values supplied or NULL if none are supplied.

Usage

```
alldiffs(predictions, differences = NULL, p.differences = NULL,
         sed = NULL, LSD = NULL, backtransforms = NULL,
         response = NULL, response.title = NULL,
         term = NULL, classify = NULL, tdf = NULL)
```

Arguments

| | |
|----------------|---|
| predictions | a data.frame containing the predicted values that is consistent with an object of class asremlPredict such as is stored in the pvals component of the prediction component of the value produced by predict.asreml. That is, in addition to variables classifying the predictions, it will include columns named predicted.value, standard.error and est.status; each row contains a single predicted value. It may also contain columns for the lower and upper confidence limits for the predictions. If LSD is not NULL, the mean LSD will be added as an attribute named meanLSD. |
| differences | a matrix containing all pairwise differences between the predictions; it should have the same number of rows and columns as there are rows in predictions. |
| p.differences | a matrix containing p-values for all pairwise differences between the predictions; each p-value is computed as the probability of a t-statistic as large as or larger than the observed difference divided by its standard error. The degrees of freedom of the t distribution for computing it are computed as the denominator degrees of freedom of the F value for the fixed term, if available; otherwise, the degrees of freedom stored in the attribute tdf are used; the matrix should be of the same size as that for differences. |
| sed | a matrix containing the standard errors of all pairwise differences between the predictions; they are used in computing the p-values. |
| LSD | a data.frame containing the mean, minimum and maximum LSD for determining the significance of pairwise differences. |
| backtransforms | a data.frame containing the backtransformed values of the predicted values that is consistent with an object of class asremlPredict such as is stored in the pvals component of the prediction component of the value produced by predict.asreml. That is, in addition to variables classifying the predictions, it will include columns named backtransformed.predictions and est.status; it may also contain columns for the lower and upper confidence limits; each row contains a single predicted value. |

| | |
|----------------|--|
| response | a character specifying the response variable for the predictions. It is stored as an attribute to the alldiffs object. |
| response.title | a character specifying the title for the response variable for the predictions. It is stored as an attribute to the alldiffs object. |
| term | a character string giving the variables that define the term that was fitted using asreml and that corresponds to classify. It is often the same as classify. It is stored as an attribute to the alldiffs object. |
| classify | a character string giving the variables that define the margins of the multiway table used in the prediction. Multiway tables are specified by forming an interaction type term from the classifying variables, that is, separating the variable names with the : operator. It is stored as an attribute to the alldiffs object. |
| tdf | an integer specifying the degrees of freedom of the standard error. It is used as the degrees of freedom for the t-distribution on which p-values and confidence intervals are based. It is stored as an attribute to the alldiffs object. |

Value

An object of S3-class alldiffs with attributes response, response.title, term, classify and tdf.

See Also

[asremlPlus-package](#), [print.alldiffs](#), [predictparallel.asreml](#), [predictionplot.asreml](#), [predictiondiffs.asreml](#), [pred.present.asreml](#)

Examples

```
## Not run:
Var.pred <- predict(current.asr, classify="Variety", sed=TRUE)$predictions
wald.tab <- current.asrt$wald.tab
den.df <- wald.tab[match("Variety", rownames(wald.tab)), "denDF"]
Var.diffs <- alldiffs(predictions = Var.pred$pvals,
                     sed = Var.pred$sed,
                     tdf = den.df)
Var.diffs <- predictiondiffs.asreml(classify = "Variety",
                                   alldiffs.obj=Var.diffs)
print.alldiffs(Var.diffs, which="differences")

## End(Not run)
```

angular

Applies the angular transformation to proportions.

Description

Applies the angular transformation to numeric values. It is given by $\sin^{-1}(\sqrt{\text{proportions}})$

Usage

```
angular(proportions, n)
```

Arguments

| | |
|-------------|------------------------------------|
| proportions | The proportions. |
| n | The divisor(s) for each proportion |

Value

A numeric.

See Also

[angular.mod](#), [power.transform](#).

Examples

```
n <-25
y <- rbinom(10, n, 0.5)
y <- c(y,0,n)
p <- y/n
p.ang <- angular(p, n)
```

`angular.mod`

Applies the modified angular transformation to a vector of counts.

Description

Applies the angular transformation to a vector of counts. A modified transformation is used that is appropriate when $N < 50$ and the proportion is not between 0.3 and 0.7. The transformation is given by $\sin^{-1} \frac{\text{count}+0.375}{n+0.75}$ `arcsin(sqrt((count + 0.375) / (n + 0.75)))`.

Usage

```
angular.mod(count, n)
```

Arguments

| | |
|-------|--|
| count | The numeric vector of counts. |
| n | The number(s) of observations from which the count(s) were obtained. |

Value

A numeric vector.

See Also

[angular](#), [power.transform](#).

Examples

```
n <-25
y <- rbinom(10, n, 0.5)
y <- c(y,0,n)
p.ang.mod <- angular.mod(y, n)
```

asremlPlus-deprecated *Deprecated Functions in Package asremlPlus*

Description

These functions have been renamed and deprecated in asremlPlus: `addrm.terms.asreml` replaced by `addrm.terms.asrtests`, `choose.model.asreml` by `addrm.terms.asrtests`, `recalc.wald.tab.asreml` by `recalc.wald.tab.asrtests`, `rmboundary.asreml` by `rmboundary.asrtests`, `sig.devn.reparam.asreml` by `sig.devn.reparam.asrtests`, `testranfix.asreml` by `testranfix.asrtests`, `testrcov.asreml` by `testrcov.asrtests`, `testswapran.asreml` by `testswapran.asrtests`, `info.crit` by `info.crit.asreml`, `reml.lrt` by `reml.lrt.asreml`

Usage

```
addrm.terms.asreml(...)
choose.model.asreml(...)
recalc.wald.tab.asreml(...)
rmboundary.asreml(...)
sig.devn.reparam.asreml(...)
testranfix.asreml(...)
testrcov.asreml(...)
testswapran.asreml(...)
info.crit(...)
reml.lrt(...)
```

Arguments

... absorbs arguments passed from the old functions of the style `foo.bar()`.

Author(s)

Chris Brien

| | |
|----------|---|
| asrtests | <i>Forms an object of S3-class asrtests that stores a fitted asreml object, a pseudo-anova table for the fixed and a history of changes and hypothesis testing used in obtaining the model.</i> |
|----------|---|

Description

An object of S3-class `asrtests` consists of a list containing:

1. `asreml.obj`: an `asreml` object containing the fit of the model;
2. `wald.tab`: a `data.frame` containing a pseudo-anova table for the fixed terms produced by `wald.asreml`, which will be called if `wald.tab` is `NULL`;

3. `test.summary`: a `data.frame` with columns `term`, `DF`, `denDF`, `p` and `action`. A row is added to it for each term that is dropped, added or tested or a note that several terms have been added or removed. A row contains the name of the term, the DF, the p-value and the action taken. Possible codes are: Dropped, Retained, Swapped, Unswapped, Significant, Nonsignificant, Absent, Added, Removed and Boundary. If the changed model did not converge, Unconverged will be added to the code. Note that the logical `asreml.obj$converge` also reflects whether there is convergence.

A call to `asrtests` with `test.summary = NULL` re-initializes the `test.summary` `data.frame`.

Usage

```
asrtests(asreml.obj, wald.tab = NULL, test.summary = NULL,
         denDF = "default", ...)
```

Arguments

| | |
|---------------------------|--|
| <code>asreml.obj</code> | an <code>asreml</code> object for a fitted model. |
| <code>wald.tab</code> | a <code>data.frame</code> containing a pseudo-anova table for the fixed terms produced by <code>wald.asreml</code> ; it should have 4 columns. Sometimes <code>wald.asreml</code> returns a <code>data.frame</code> and at other times a list. For example, it may return a list when <code>denDF</code> is used. In this case, the Wald component of the list is to be extracted and stored. |
| <code>test.summary</code> | a <code>data.frame</code> with columns <code>term</code> , <code>DF</code> , <code>denDF</code> , <code>p</code> and <code>action</code> containing the results of previous hypothesis tests. |
| <code>denDF</code> | Specifies the method to use in computing approximate denominator degrees of freedom when <code>wald.asreml</code> is called. Can be <code>none</code> to suppress the computations, <code>numeric</code> for numerical methods, <code>algebraic</code> for algebraic methods or <code>default</code> , the default, to automatically choose numeric or algebraic computations depending on problem size. The denominator degrees of freedom are calculated according to Kenward and Roger (1997) for fixed terms in the dense part of the model. |
| <code>...</code> | further arguments passed to <code>wald.asreml</code> . |

Value

An object of S3-class `asrtests`.

See Also

[asremlPlus-package](#), [testranfix.asrtests](#), [choose.model.asrtests](#), [rmboundary.asrtests](#), [sig.devn.reparam.asrtests](#)

Examples

```
## Not run:
data(Wheat.dat)

# Fit initial model
current.asr <- asreml(yield ~ Rep + WithinColPairs + Variety,
                     random = ~ Row + Column + units,
                     rcov = ~ ar1(Row):ar1(Column),
```

```

data=Wheat.dat)

# Load current fit into an asrtests object
current.asrt <- asrtests(current.asr, NULL, NULL)

# Check for and remove any boundary terms
current.asrt <- rmboundary.asrtests(current.asrt)

## End(Not run)

```

`choose.model.asrtests` *Determines the set of significant terms taking into account the hierarchy or marginality relations.*

Description

Performs a series of hypothesis tests taking into account the marginality of terms. In particular, a term will not be tested if it is marginal to (or nested in) one that is significant. For example, if A:B is significant, then neither A nor B will be tested. For a random term, the term is removed from the model fit, any boundary terms are removed using `rmboundary.asrtests` and a REML likelihood ratio test is performed using `reml.lrt.asreml`. If it is not significant and `drop.ran.ns` is TRUE, the term is permanently removed from the model. Note that if boundary terms are removed, the reduced model may not be nested in the full model in which case the test is not valid. For fixed terms, the Wald tests are performed and the p-value for the term obtained. If it is not significant and `drop.fix.ns` is TRUE, the term is permanently removed from the model. A row is added to `test.summary` for each term that is tested.

Usage

```

choose.model.asrtests(terms.marginality=NULL, asrtests.obj,
  alpha = 0.05, drop.ran.ns=TRUE, positive.zero = FALSE,
  drop.fix.ns=FALSE, denDF = "default", dDF.na = "none",
  dDF.values = NULL, trace = FALSE, update = TRUE,
  set.terms = NULL, ignore.suffices = TRUE,
  constraints = "P", initial.values = NA, ...)

```

Arguments

| | |
|--------------------------------|--|
| <code>terms.marginality</code> | a square matrix of ones and zeros with row and column names being the names of the terms. The diagonal elements should be one, indicating that a term is marginal to itself. Elements should be one if the row term is marginal to the column term. All other elements should be zero. |
| <code>asrtests.obj</code> | an <code>asrtests</code> object for a fitted model that is a list containing an <code>asreml</code> object, a <code>wald.tab</code> data.frame with 4 columns, and a <code>data.frame</code> with 5 columns that records any previous changes and tests in the fitted model. |
| <code>alpha</code> | the significance level for the test. |
| <code>drop.ran.ns</code> | a logical indicating whether to drop nonsignificant random terms from the model. |
| <code>positive.zero</code> | a logical indicating whether the hypothesized values for the variance components being tested are on the boundary of the parameter space. For example, this is be true for positively-constrained variance components that, under the reduced model, are zero. |

| | |
|-----------------|--|
| drop.fix.ns | a logical indicating whether to drop a fixed term from the model when it is nonsignificant |
| denDF | Specifies the method to use in computing approximate denominator degrees of freedom when <code>wald.asreml</code> is called. Can be <code>none</code> to suppress the computations, <code>numeric</code> for numerical methods, <code>algebraic</code> for algebraic methods or <code>default</code> , the default, to automatically choose numeric or algebraic computations depending on problem size. The denominator degrees of freedom are calculated according to Kenward and Roger (1997) for fixed terms in the dense part of the model. |
| dDF.na | the method to use to obtain substitute denominator degrees of freedom. when the numeric or algebraic methods produce an NA. If <code>dDF.na = "none"</code> , no substitute denominator degrees of freedom are employed; if <code>dDF.na = "residual"</code> , the residual degrees of freedom from <code>asreml.obj\$nedf</code> are used; if <code>dDF.na = "maximum"</code> , the maximum of those <code>denDF</code> that are available, excluding that for the Intercept, is used; if all <code>denDF</code> are NA, <code>asreml.obj\$nedf</code> is used. If <code>dDF.na = "supplied"</code> , a vector of values for the denominator degrees of freedom is to be supplied in <code>dDF.values</code> . Any other setting is ignored and a warning message produced. Generally, substituting these degrees of freedom is anticonservative in that it is likely that the degrees of freedom used will be too large. |
| dDF.values | A vector of values to be used when <code>dDF.na = "supplied"</code> . Its values will be used when <code>denDF</code> in a test for a fixed effect is NA. This vector must be the same length as the number of fixed terms, including (Intercept) whose value could be NA. |
| trace | if TRUE then partial iteration details are displayed when <code>ASReml-R</code> functions are invoked; if FALSE then no output is displayed. |
| update | if TRUE then <code>update.asreml</code> is called in testing models. In doing this the arguments <code>R.param</code> and <code>G.param</code> are set to those in the <code>asreml</code> object stored in <code>asrtests.obj</code> so that the values from the previous model are used as starting values. If FALSE then a call is made to <code>asreml</code> in which the only changes to the <code>asreml.obj</code> stored in the supplied <code>asrtests.obj</code> are (i) to the terms in the fixed and random models corresponding to terms in <code>terms.marginality</code> and (ii) those modifications specified via <code>...</code> |
| set.terms | a character vector specifying the terms that are to have constraints and/or initial values set prior to fitting. |
| ignore.suffices | a logical vector specifying whether the suffices of the <code>asreml</code> -assigned names of the variance terms (i.e. the information to the right of an "!", other than "R!") is to be ignored in matching elements of terms. If TRUE for an element of terms, the suffices are stripped from the <code>asreml</code> -assigned names. If FALSE for an element of terms, the element must exactly match an <code>asreml</code> -assigned name for a variance term. This vector must be of length one or the same length as <code>terms</code> . If it is of length one then the same action is applied to the <code>asreml</code> -assigned suffices for all the terms in <code>terms</code> . |
| constraints | a character vector specifying the constraints to be applied to the terms specified in <code>terms</code> . This vector must be of length one or the same length as <code>terms</code> . If it is of length one then the same constraint is applied to all the terms in <code>terms</code> . If any of the constraints are equal to NA then they are left unchanged for those terms. |
| initial.values | a character vector specifying the initial values for the terms specified in <code>terms</code> . This vector must be of length one or the same length as <code>terms</code> . If it is of length one then the same initial value is applied to all the terms in <code>terms</code> . If any of the <code>initial.values</code> are equal to NA then they are left unchanged for those terms. |

... further arguments passed to `asreml` and `wald.asreml` via `testranfix.asrtests`.

Value

A list containing:

1. `asrtests.obj`: an `asrtests` object, containing the `asreml` object corresponding to the final fit, a `wald.tab` data.frame, and a `test.summary` data.frame that contains a record of the testing of the terms (see `asrtests` for more details);
2. `sig.tests`: a character vector whose elements are the the significant terms amongst those tested.

See Also

`asrtests`, `testranfix.asrtests`, `testrcov.asrtests`, `reml.lrt.asreml`, `rmboundary.asrtests`, `newfit.asreml`, `addrm.terms.asrtests`, `sig.devn.reparam.asrtests`

Examples

```
terms.treat <- c("Sources", "Type", "Species",
                "Sources:Type", "Sources:Species")
terms <- sapply(terms.treat,
               FUN=function(term){paste("spl(xDay):", term, sep="")},
               simplify=TRUE)
terms <- c("spl(xDay)", terms)
marginality <- matrix(c(1,0,0,0,0,0, 1,1,0,0,0,0, 1,0,1,0,0,0,
                        1,0,1,1,0,0, 1,1,1,1,1,0, 1,1,1,1,1,1), nrow=6)
rownames(marginality) <- terms
colnames(marginality) <- terms
## Not run: choose <- choose.model.asrtests(marginality, current.asrt,
                                           pos=TRUE)
current.asrt <- choose$asrtests.obj

## End(Not run)
```

info.crit.asreml

Computes AIC and BIC for a model.

Description

Computes Akiake and Bayesian (Schwarz) Information Criteria for a model. The function `info.crit` is provided for backwards compatibility.

Usage

```
info.crit.asreml(asreml.obj)
```

Arguments

`asreml.obj` An `asreml` object resulting from the fitting of a model using REML.

Value

A data frame containing the Residual degrees of freedom, AIC, BIC and log of the REML value.

See Also[reml.lrt.asreml](#)**Examples**

```
## Not run:
data(Wheat.dat)
current.asr <- asreml(yield ~ Rep + WithinColPairs + Variety,
  random = ~ Row + Column + units,
  rcov = ~ ar1(Row):ar1(Column),
  data=Wheat.dat)
info.crit.asreml(current.asr)

## End(Not run)
```

| | |
|---------------|--|
| newfit.asreml | <i>refits an asreml model with modified model formula using either a call to update.asreml or a direct call to asreml.</i> |
|---------------|--|

Description

Extracts the call from the `asreml.obj` and evaluates that call, replacing any arguments with changed values. If `update` is `TRUE` and `set.terms` is not set, the call is evaluated using `update.asreml`; otherwise, it is evaluated using a direct call to `asreml`. The principal difference is that the latter does not enforce the use of previous values of the variance parameters as initial values; it sets `G.param` and `R.param` to `NULL` or to values as specified for `set.terms`. The `...` argument can be used to pass `G.param` and/or `R.param`, provided `update` is `FALSE` and `set.terms` is not set.

Usage

```
newfit.asreml(asreml.obj, fixed., random., sparse., rcov.,
  update = TRUE, keep.order = TRUE, set.terms = NULL,
  ignore.suffices = TRUE, constraints = "P", initial.values = NA, ...)
```

Arguments

| | |
|-------------------------|---|
| <code>asreml.obj</code> | a valid <code>asreml</code> object with with a component named <code>call</code> (from a previous call to either <code>asreml</code> or <code>update.asreml</code>). |
| <code>fixed.</code> | a character or formula specifying changes to the fixed formula. This is a two-sided formula where <code>"."</code> is substituted for existing components in the fixed component of <code>asreml.obj\$call</code> . |
| <code>random.</code> | a character or formula specifying changes to the random formula. This is a one-sided formula where <code>"."</code> is substituted for existing components in the random component of <code>asreml.obj\$call</code> . |
| <code>sparse.</code> | a character or formula specifying changes to the sparse formula. This is a one-sided formula where <code>"."</code> is substituted for existing components in the sparse component of <code>asreml.obj\$call</code> . |
| <code>rcov.</code> | a character or formula specifying changes to the error formula. This is a one-sided formula where <code>"."</code> is substituted for existing components in the <code>rcov</code> component of <code>asreml.obj\$call</code> . |

| | |
|-----------------|---|
| update | a logical indicated whether to use <code>update.asreml</code> or <code>asreml</code> to evaluate the modified call. If TRUE, use <code>update.asreml</code> to evaluate the modified call. In doing this the arguments <code>R.param</code> and <code>G.param</code> are set to those in the <code>asreml.obj</code> so that the values from the previous model are used as starting values. If FALSE then a call is made to <code>asreml</code> itself, in which the only changes from the previous call are those specified in the arguments to <code>newfit.asreml</code> . |
| keep.order | a logical value indicating whether the terms should keep their positions. If FALSE the terms are reordered so that main effects come first, followed by the interactions, all second-order, all third-order and so on. Effects of a given order are kept in the order specified. |
| set.terms | a character vector specifying the terms that are to have constraints and/or initial values set prior to fitting. |
| ignore.suffices | a logical vector specifying whether the suffices of the <code>asreml</code> -assigned names of the variance terms (i.e. the information to the right of an "!", other than "R!") is to be ignored in matching elements of terms. If TRUE for an element of terms, the suffices are stripped from the <code>asreml</code> -assigned names. If FALSE for an element of terms, the element must exactly match an <code>asreml</code> -assigned name for a variance term. This vector must be of length one or the same length as <code>terms</code> . If it is of length one then the same action is applied to the <code>asreml</code> -assigned suffices for all the terms in <code>terms</code> . |
| constraints | a character vector specifying the constraints to be applied to the terms specified in <code>terms</code> . This vector must be of length one or the same length as <code>terms</code> . If it is of length one then the same constraint is applied to all the terms in <code>terms</code> . If any of the constraints are equal to NA then they are left unchanged for those terms. |
| initial.values | a character vector specifying the initial values for the terms specified in <code>terms</code> . This vector must be of length one or the same length as <code>terms</code> . If it is of length one then the same initial value is applied to all the terms in <code>terms</code> . If any of the <code>initial.values</code> are equal to NA then they are left unchanged for those terms. |
| ... | additional arguments to the call, or arguments with changed values. |

Value

An `asreml` object.

References

Butler, D. G., et al. (2010). *Analysis of Mixed Models for S language environments: ASReml-R reference manual*. Brisbane, DPI Publications.

See Also

`update.asreml`, [setvarianceterms.asreml](#)

Examples

```
## Not run:
m2.asreml <- newfit.asreml(m1.asreml, random. = "~ . - Blocks:Plots", maxiter=75)

## End(Not run)
```

| | |
|------------------|--|
| newrcov.asrtests | <i>Fits a new rcov formula using asreml.</i> |
|------------------|--|

Description

Fits a new rcov formula using asreml. Any boundary terms resulting after fitting the new rcov model are removed using [rmboundary.asrtests](#). A row is added to the test.summary data.frame using the supplied label and the action Set.

Usage

```
newrcov.asrtests(terms=NULL, asrtests.obj, label = "R model",
                 update = TRUE, trace = FALSE, denDF="default",
                 set.terms = NULL, ignore.suffices = TRUE,
                 constraints = "P", initial.values = NA, ...)
```

Arguments

| | |
|-----------------|---|
| terms | a model for the rcov argument in asreml, stored as a character. |
| asrtests.obj | an asrtests object for a fitted model that is a list containing an asreml object, a wald.tab data.frame with 4 columns, and a data.frame with 5 columns that records any previous changes and tests in the fitted model. |
| label | a character string to use as the label in test.summary and which indicates what is being tested. |
| update | if TRUE then update.asreml is called to fit the model with the rcov model supplied in terms. In doing this the arguments R.param and G.param are set to those in the asreml object stored in asrtests.obj so that the values from the previous model are used as starting values. If FALSE then a call is made to asreml in which the only changes from the previous call are that (i) rcov model is that specified in terms and (ii) modifications specified via ... are made. |
| trace | if TRUE then partial iteration details are displayed when ASReML-R functions are invoked; if FALSE then no output is displayed. |
| denDF | Specifies the method to use in computing approximate denominator degrees of freedom when wald.asreml is called. Can be none to suppress the computations, numeric for numerical methods, algebraic for algebraic methods or default, the default, to automatically choose numeric or algebraic computations depending on problem size. The denominator degrees of freedom are calculated according to Kenward and Roger (1997) for fixed terms in the dense part of the model. |
| set.terms | a character vector specifying the terms that are to have constraints and/or initial values set prior to fitting. |
| ignore.suffices | a logical vector specifying whether the suffices of the asreml-assigned names of the variance terms (i.e. the information to the right of an "!", other than "R!") is to be ignored in matching elements of terms. If TRUE for an element of terms, the suffices are stripped from the asreml-assigned names. If FALSE for an element of terms, the element must exactly match an asreml-assigned name for a variance term. This vector must be of length one or the same length as terms. If it is of length one then the same action is applied to the asreml-assigned suffices for all the terms in terms. |

| | |
|----------------|---|
| constraints | a character vector specifying the constraints to be applied to the terms specified in terms. This vector must be of length one or the same length as terms. If it is of length one then the same constraint is applied to all the terms in terms. If any of the constraints are equal to NA then they are left unchanged for those terms. |
| initial.values | a character vector specifying the initial values for the terms specified in terms. This vector must be of length one or the same length as terms. If it is of length one then the same initial value is applied to all the terms in terms. If any of the initial.values are equal to NA then they are left unchanged for those terms. |
| ... | further arguments passed to asreml and to wald.asreml. |

Value

An `asrtests` object, which is a list containing:

1. `asreml.obj`: an `asreml` object containing the fit after the term has been omitted from the model;
2. `wald.tab`: a 4-column data.frame containing a pseudo-anova table for the fixed terms produced by `wald.asreml`;
3. `test.summary`: a data.frame with columns term, DF, denDF, p and action. A row is added to it for each term that is dropped, added or tested or a note that several terms have been added or removed. A row contains the name of the term, the DF, the p-value and the action taken. Possible codes are: Dropped, Retained, Swapped, Unswapped, Significant, Nonsignificant, Absent, Added, Removed and Boundary. If the changed model did not converge, Unconverged will be added to the code. Note that the logical `asreml.obj$converge` also reflects whether there is convergence.

If the term is not in the model, then the supplied `asreml` object will be returned. Also, `reml.test` will have the likelihood ratio and the p-value set to NA and the degrees of freedom to zero. Similarly, the row of `test.summary` for the term will have its name, a p-value set to NA, and action set to Absent.

See Also

[asremlPlus-package](#), [asrtests](#), [testrcov.asrtests](#),
[choose.model.asrtests](#), [reml.lrt.asreml](#), [rmboundary.asrtests](#),
[newfit.asreml](#), [testswapran.asrtests](#), [addrm.terms.asrtests](#),
[sig.devn.reparam.asrtests](#)

Examples

```
## Not run:
data(Wheat.dat)
current.asr <- asreml(yield ~ Rep + WithinColPairs + Variety,
                     random = ~ Row + Column + units,
                     rcov = ~ ar1(Row):ar1(Column),
                     data=Wheat.dat)
current.asrt <- asrtests(current.asr, NULL, NULL)
current.asrt <- rmboundary.asrtests(current.asrt)
# Fit Row autocorrelation
current.asrt <- newrcov.asrtests("~ Row:ar1(Column)", current.asrt,
                                label="Row autocorrelation")
print(current.asrt)

## End(Not run)
```

| | |
|------------|--|
| num.recode | <i>Recodes the unique values of a vector using the values in a new vector.</i> |
|------------|--|

Description

Recodes the unique values of a variate using the value in position *i* of the `new.values` vector to replace the *i*th sorted unique values of `x`. The new levels do not have to be unique.

Usage

```
num.recode(x, new.values)
```

Arguments

| | |
|-------------------------|---|
| <code>x</code> | The vector to be recoded. |
| <code>new.values</code> | A vector of length <code>unique(x)</code> containing values to use in the recoding. |

Value

A vector.

Author(s)

Chris Brien

See Also

`dae::fac.recode`.

Examples

```
## set up a factor with labels
x <- rep(c(-42, -14, 14, 42), 4)

## recode x
b <- num.recode(x, c(0, 28, 56, 84))
```

| | |
|----------------|--|
| permute.square | <i>Permutes the rows and columns of a square matrix.</i> |
|----------------|--|

Description

Permutes the rows and columns of a square matrix.

Usage

```
permute.square(x, permutation)
```

Arguments

`x` A square matrix.
`permutation` A vector specifying the new order of rows and columns.

Value

A square matrix.

See Also

[permute.to.zero.lowertri](#)

Examples

```
terms.marginality <- matrix(c(1,0,0,0,0, 0,1,0,0,0, 0,1,1,0,0,
                             1,1,1,1,0, 1,1,1,1,1), nrow=5)
permtn <- c(1,3,2,4,5)
terms.marginality <- permute.square(terms.marginality, permtn)
```

`permute.to.zero.lowertri`

Permutes a square matrix until all the lower triangular elements are zero.

Description

Permutes a square matrix until all the lower triangular elements are zero.

Usage

```
permute.to.zero.lowertri(x)
```

Arguments

`x` A square matrix of order n with at least $n*(n-1)/2$ zero elements.

Value

A square matrix.

See Also

[permute.square](#)

Examples

```
terms.marginality <- matrix(c(1,0,0,0,0, 0,1,0,0,0, 0,1,1,0,0,
                             1,1,1,1,0, 1,1,1,1,1), nrow=5)
terms.marginality <- permute.to.zero.lowertri(terms.marginality)
```

plotvariofaces.asreml *plot empirical variogram faces, including envelopes, from supplied residuals as described by Stefanova, Smith & Cullis (2009)*

Description

Produces a plot for each face of an empirical 2D variogram based on supplied residuals from both an observed data set and simulated data sets. Those from simulated data sets are used to produce confidence envelopes. If the data consists of sections, such as separate experiments, the two variogram faces are produced for each section. This function is less efficient in storage terms than `variofaces.asreml`, because here the residuals from all simulated data sets must be saved, in addition to the values for the variogram faces; in `variofaces.asreml`, the residuals for each simulated data set are discarded after the variogram has been calculated. On the other hand, the present function is more flexible, because there is no restriction on how the residuals are obtained.

Usage

```
plotvariofaces.asreml(data, residuals, restype="Residuals", ...)
```

Arguments

| | |
|-----------|--|
| data | A data.frame with either 3 or 4 columns. Only if there are 4 columns, the first should be a factor indexing sections for which separate variogram plots are to be produced. In either case, the other 3 columns should be, in order, (i) a factor indexing the x-direction, (ii) a factor indexing the y-direction, and (iii) the residuals for the observed response. |
| residuals | A data.frame, with either 2 or 3 initial columns followed by columns, each of which are the residuals from a simulated data set. |
| restype | A character describing the type of residuals that have been supplied. It will be used in the plot titles. |
| ... | Other arguments that are passed down to the function <code>asreml.variogram</code> . |

Details

For each set of residuals, `asreml.variogram` is used to obtain the empirical variogram, from which the values for its faces are obtained. Plots are produced for each face and include the observed residuals and the 2.5%, 50% & 97.5% quantiles.

Value

A list with the following components:

1. **face1**: a data.frame containing the variogram values on which the plot for the first dimension is based.
2. **face2**: a data.frame containing the variogram values on which the plot for the second dimension is based.

Author(s)

Chris Brien

References

Stefanova, K. T., Smith, A. B. & Cullis, B. R. (2009) Enhanced diagnostics for the spatial analysis of field trials. *Journal of Agricultural, Biological, and Environmental Statistics*, **14**, 392–410.

See Also

[asremlPlus-package](#), [asreml](#), [asreml.variogram](#), [variofaces.asreml](#), [simulate.asreml](#).

Examples

```
## Not run:
data(Wheat.dat)
current.asr <- asreml(yield ~ Rep + WithinColPairs + Variety,
                     random = ~ Row + Column + units,
                     rcov = ~ ar1(Row):ar1(Column),
                     data=Wheat.dat)
current.asrt <- asrtests(current.asr, NULL, NULL)
current.asrt <- rmboundary.asrtests(current.asrt)
# Form variance matrix based on estimated variance parameters
s2 <- current.asr$sigma2
gamma.Row <- current.asr$gammas[1]
gamma.unit <- current.asr$gammas[2]
rho.r <- current.asr$gammas[4]
rho.c <- current.asr$gammas[5]
row.ar1 <- mat.ar1(order=10, rho=rho.r)
col.ar1 <- mat.ar1(order=15, rho=rho.c)
V <- gamma.Row * fac.sumop(Wheat.dat$Row) +
    gamma.unit * diag(1, nrow=150, ncol=150) +
    mat.dirprod(col.ar1, row.ar1)
V <- s2*V

#Produce variogram faces plot (Stefanova et al, 2009)
resid <- simulate.asreml(current.asr, V=V, which="residuals")
resid$residuals <- cbind(resid$observed[c("Row", "Column")],
                        resid$residuals)
plotvariofaces.asreml(data=resid$observed[c("Row", "Column", "residuals")],
                      residuals=resid$residuals,
                      restype="Standardized conditional residuals")

## End(Not run)
```

power.transform

Perform a combination of a linear and power transformation on a variable whose name is given as a character string in var.name. The transformed variable is stored in the data.frame data.

Description

Perform a combination of a linear and a power transformation on a variable whose name is given as a character string in var.name. The transformed variable is stored in the data.frame data. The name of the transformed variable is made by prepending to the original var.name a combination of (i) .offset, if offset is nonzero, (ii) neg., if scale is -1, or scaled., if abs(scale) is other than one, and (iii) either log., sqrt., recip. or power., if power is other than one. No action

is taken if there is no transformation (i.e. `offset = 0`, `scale = 1` and `power = 1`). Also, the `titles` list is extended to include a component with a generated title for the transformed variable with text indicating the transformation prepended to the title for the `var.name` obtained from the `titles` list. For nonzero `offset`, 'Offset ' is prepended, For scaled not equal to one, the possible prepends are 'Negative of ' and 'Scaled '. The possible prepended texts for power not equal to one are 'Logarithm of', 'Square root of', 'Reciprocal of' and 'Power nnnn of', where `nnn` is the power used.

Usage

```
power.transform(var.name, power = 1, offset = 0, scale = 1, titles = NULL, data)
```

Arguments

| | |
|-----------------------|---|
| <code>var.name</code> | A character string specifying the name of the variable in the <code>data.frame</code> <code>data</code> that is to be transformed. |
| <code>power</code> | A number specifying the power to be used in the transformation. If equal to 1, the default, no power transformation is applied. Otherwise, the variable is raised to the specified power, after scaling and applying any nonzero <code>offset</code> . If <code>power = 0</code> , the natural logarithm is used to transform the response; however, if the smallest value to be log-transformed is less than 1e-04, an error is generated. A log-transformation in this situation may be possible if a nonzero <code>offset</code> and/or a <code>scale</code> not equal to one is used. |
| <code>offset</code> | A number to be added to each value of the variable, after any scaling and before applying any power transformation. |
| <code>scale</code> | A number to multiply each value of the variable, before adding any <code>offset</code> and applying any power transformation. |
| <code>titles</code> | A character vector, each element of which is named for a variable in <code>data</code> and is a character string giving a title to use in output (e.g. tables and graphs) involving the variable. If titles are not supplied, the column name of the variable in <code>dta</code> is used. |
| <code>data</code> | A <code>data.frame</code> containing the variable to be transformed and to which the transformed variable is to be appended. |

Value

A list with a component named `data` that is the `data.frame` containing the transformed variable, a component named `tvar.name` that is a character string that is the name of the transformed variable in `data`, and a component named `titles` that extends the list supplied in the `titles` argument to include a generated title for the transformed title, the name of the new component being `tvar.name`.

Author(s)

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See Also

[angular](#), [angular.mod](#).

Examples

```
## set up a factor with labels
x.dat <- data.frame(y = c(14, 42, 120, 150))

## transform y to logarithms
trans <- power.transform("y", power = 0, titles=list(y = "Length (cm)"), data = x.dat)
x.dat <- trans$data
tvar.name <- trans$tvar.name

## transform y to logarithms after multiplying by -1 and adding 1.
z.dat <- data.frame( y = c(-5.25, -4.29, -1.22, 0.05))
trans <- power.transform("y", power = 0, scale = -1, offset = 1 ,
                        titles=list(y = "Potential"), data = z.dat)

z.dat <- trans$data
tvar.name <- trans$tvar.name
```

| | |
|---------------------|---|
| pred.present.asreml | <i>This function forms the predictions for each significant term and presents them in tables and/or graphs.</i> |
|---------------------|---|

Description

This function forms the predictions for each term in terms using a supplied asreml object and [predictparallel.asreml](#). Tables are produced using [predictparallel.asreml](#), in conjunction with [predictiondiffs.asreml](#), with the argument tables specifying which tables are printed. The argument plots, along with transform.power, controls which plots are produced. The plots are produced using [predictionplot.asreml](#), with line plots produced when variables involving x.num or x.fac are involved in classify for the predictions and bar charts otherwise. In order to get the correct predictions you may need to supply additional arguments to predict through ... e.g. present.

Usage

```
pred.present.asreml(terms, asreml.obj = NULL,
                    wald.tab = NULL, dDF.na = "residual", dDF.values = NULL,
                    x.num = NULL, x.fac = NULL, nonx.fac.order = NULL,
                    x.pred.values = NULL, x.plot.values = NULL,
                    plots = "predictions", panels = "multiple",
                    graphics.device = NULL,
                    error.intervals = "Confidence", avsed.tolerance = 0.25,
                    titles = NULL, colour.scheme = "colour", save.plots = FALSE,
                    transform.power = 1, offset = 0, scale = 1,
                    pairwise = TRUE, tables = "all", levels.length = NA,
                    alpha = 0.05, inestimable.rm = TRUE, trace = FALSE, ...)
```

Arguments

| | |
|------------|---|
| terms | a character vector giving the terms for which predictions are required. |
| asreml.obj | asreml object for a fitted model. |
| wald.tab | a data frame containing the pseudo-anova table for the fixed terms produced by a call to wald.asreml. The main use of it here is in getting denominator degrees of freedom when confidence intervals are to be plotted. |

| | |
|-----------------|---|
| dDF.na | the method to use to obtain approximate denominator degrees of freedom. when the numeric or algebraic methods produce an NA. Consistent with when no denDF are available, the default is "residual" and so the residual degrees of freedom from asreml.obj\$nedf are used. If dDF.na = "none", no substitute denominator degrees of freedom are employed; if dDF.na = "maximum", the maximum of those denDF that are available, excluding that for the Intercept, is used; if all denDF are NA, asreml.obj\$nedf is used. If dDF.na = "supplied", a vector of values for the denominator degrees of freedom is to be supplied in dDF.values. Any other setting is ignored and a warning message produced. Generally, substituting these degrees of freedom is anticonservative in that it is likely that the degrees of freedom used will be too large. |
| dDF.values | A vector of values to be used when dDF.na = "supplied". Its values will be used when denDF in a test for a fixed effect is NA. This vector must be the same length as the number of fixed terms, including (Intercept) whose value could be NA. |
| x.num | A character string giving the name of the numeric covariate that corresponds to x.fac, is potentially included in terms in the fitted model and which corresponds to the x-axis variable. It should have the same number of unique values as the number of levels in x.fac. |
| x.fac | A character string giving the name of the factor that corresponds to x.num, is potentially included in terms in the fitted model and which corresponds to the x-axis variable. It should have the same number of levels as the number of unique values in x.num. The levels of x.fac must be in the order in which they are to be plotted - if they are dates, then they should be in the form yyyyymmdd which can be achieved using as.Date. |
| nonx.fac.order | A character vector giving the order in which factors other than x.fac are to be plotted in plots with multiple panels (i.e. where the number of non-x factors is greater than 1). The first factor in the vector will be plotted on the X axis (if there is no x.num or x.fac. Otherwise, the order of plotting the factors is in columns (X facets) and then rows (Y facets). By default the order is in decreasing order for the numbers of levels of the non x factors. |
| x.pred.values | The values of x.num for which predicted values are required. |
| x.plot.values | The actual values to be plotted on the x axis or in the labels of tables. They are needed when values different to those in x.num are to be plotted or x.fac is to be plotted because there is no x.num term corresponding to the same term with x.fac. |
| plots | Possible values are "none", "predictions", "backtransforms" and "both". Plots are not produced if the value is "none". If data are not transformed for analysis (transform.power = 1), a plot of the predictions is produced provided plots is not "none". If the data are transformed, the value of plots determines what is produced. |
| panels | Possible values are "single" and "multiple". When line plots are to be produced, because variables involving x.num or x.fac are involved in classify for the predictions, panels determines whether or not a single panel or multiple panels in a single window are produced. The panels argument is ignored for bar charts. |
| graphics.device | A character specifying a graphics device for plotting. The default is graphics.device = NULL, which will result in plots being produced on the current graphics device. Setting it to "windows", for example, will result in a windows graphics device being opened. |

`error.intervals`

A character string indicating the type of error interval, if any, to calculate and plot in order to indicate uncertainty in the results. Possible values are "none", "StandardError", "Confidence" and "halfLeastSignificant". The default is for confidence limits to be used. The "halfLeastSignificant" option results in half the Least Significant Difference (LSD) being added and subtracted to the predictions, the LSD being calculated using the average of the standard errors of all pairwise differences (SEDs) between the predictions. However, if the range of the SEDs divided by the average of the SEDs exceeds `avsed.tolerance`, calculations and plotting will revert to confidence intervals. Also, half LSDs cannot be used for backtransformed values and so confidence intervals will be used instead.

`avsed.tolerance`

The values of the range of the SEDs divided by the average of the SEDs that, if exceeded, will cause calculations and plotting to revert to confidence intervals. It should be a value between 0 and 1.

`titles`

A list, each component of which is named for an object name and contains a character string giving a title to use in output (e.g. tables and graphs) for the object. Here they will be used for axis labels.

`colour.scheme`

A character string specifying the colour scheme for the plots. The default is "colour" which produces coloured lines and bars, a grey background and white gridlines. A value of "black" results in black lines, grey bars and gridlines and a white background.

`save.plots`

A logical that determines whether any plots will be saved. If they are to be saved, a file name will be generated that consists of the following elements separated by full stops: the response variable name with `.back` if backtransformed values are being plotted, the classify term, Bar or Line and, if `error.intervals` is not "none", one of SE, CI or LSI. The file will be saved as a 'png' file in the current work directory.

`transform.power`

A number specifying the power of a transformation, if one has been applied to the response variable. Unless it is equal to 1, the default, back-transforms of the predictions will be obtained and presented in tables or graphs as appropriate. The back-transform will raise the predictions to the power equal to the reciprocal of `transform.power`, unless it equals 0 in which case the exponential will be taken. Any scaling and offsetting will also be taken into account in the backtransformation.

`offset`

A number that has been added to each value of the response after any scaling and before applying any power transformation. Unless it is equal to 0, the default, back-transforms of the predictions will be obtained and presented in tables or graphs as appropriate. The back-transform will, after backtransforming for any power transformation and scaling for any scale transformation, subtract the offset.

`scale`

A number by which each value of the response has been multiply before adding any offset and applying any power transformation. Unless it is equal to 1, the default, back-transforms of the predictions will be obtained and presented in tables or graphs as appropriate. The back-transform will, after backtransforming for any power transformation, divide by the scale.

`pairwise`

A logical indicating whether all pairwise differences of the predictions and their standard errors and p-values are to be computed and stored. If `tables` is equal to

| | |
|----------------|---|
| | "differences" or "all" or error.intervals is equal to "halfLeastSignificant", they will be stored irrespective of the value of pairwise. |
| tables | The elements of the <code>alldiffs</code> object to print. Possible values are "none", "predictions", "backtransforms", "nodifferences", "differences" and "all". The option "nodifferences" will result in "predictions" and "backtransforms" being printed. The option "differences" will also result in the printing of the predictions from the <code>alldiffs</code> object. |
| levels.length | The maximum number of characters from the the levels of factors to use in the row and column labels of the tables produced by <code>predictiondiffs.asreml</code> . |
| alpha | the significance level for the test or 1 - alpha is the confidence level for confidence intervals. |
| inestimable.rm | A logical indicating whether rows for predictions that are not estimable are to be removed from the components of the <code>alldiffs</code> object. |
| trace | if TRUE then partial iteration details are displayed when ASReml-R functions are invoked; if FALSE then no output is displayed. |
| ... | further arguments passed to <code>predict.asreml</code> via <code>predictparallel.asreml</code> and to <code>ggplot</code> via <code>predictionplot.asreml</code> . |

Value

a list containing a `alldiffs` object for each term for which tables are produced. The names of the components of this list are the terms with full-stops (.) replacing colons (:).

See Also

`predictparallel.asreml`, `predictiondiffs.asreml`, `predictionplot.asreml`, `print.alldiffs`, `as.Date`, `Devices`

Examples

```
## Not run:
pred.present.asreml(choose$sig.terms, current.asrt$asreml.obj, current.asrt$wald.tab,
  x.fac = "Date", x.num = "xDay",
  x.pred.values=sort(unique(Runoff.longi.dat$xDay)),
  x.plot.values=c(0,28,56,84),
  x.title = "Days since first observation",
  y.title = "Predicted log(Turbidity)",
  present = c("Type","Species","Sources"))

## End(Not run)
```

predictiondiffs.asreml

Uses information in a supplied alldiffs object to forms all pairwise differences between a set of predictions, the p-values for a test of whether the differences are significantly different from zero, and the minimum, mean and maximum LSD values, provided they are not present in the supplied alldiffs object.

Description

Uses predictions and standard errors of pairwise differences from an `alldiffs` object to form, for those components not already present, (i) a table of all pairwise differences of the predictions in an `alldiffs` object, (ii) the p-values of each pairwise difference, and (iii) the minimum, mean and maximum LSD values. Predictions that are aliased (or nonestimable) are removed from the predictions component of the `alldiffs` object and standard errors of differences involving them are removed from the `sed` component.

Each p-value is computed as the probability of a t-statistic as large as or larger than the absolute value of the observed difference divided by its standard error. The p-values are stored in the `p.differences` component. The degrees of freedom of the t-distribution is the degrees of freedom stored in the `tdf` attribute of the `alldiffs` object. This t-distribution is also used in calculating the LSD statistics stored in the `alldiffs` object.

Usage

```
predictiondiffs.asreml(classify, alldiffs.obj,
                      x.num = NULL, x.fac = NULL,
                      levels.length = NA,
                      pairwise = TRUE, alpha = 0.05,
                      inestimable.rm = TRUE)
```

Arguments

| | |
|-----------------------------|--|
| <code>classify</code> | a character string giving the variables that define the margins of the multiway table to be predicted. Multiway tables are specified by forming an interaction type term from the classifying variables, that is, separating the variable names with the <code>:</code> operator. |
| <code>alldiffs.obj</code> | An <code>alldiffs</code> object for a fitted model. Note that the attribute <code>tdf</code> , being the degrees of freedom for the critical t-value to be used in computing p-values, should be set to an appropriate value. |
| <code>x.num</code> | A character string giving the name of the numeric covariate that corresponds to <code>x.fac</code> , is potentially included in terms in the fitted model and which corresponds to the x-axis variable. It should have the same number of unique values as the number of levels in <code>x.fac</code> . |
| <code>x.fac</code> | A character string giving the name of the factor that corresponds to <code>x.num</code> , is potentially included in terms in the fitted model and which corresponds to the x-axis variable. It should have the same number of levels as the number of unique values in <code>x.num</code> . The levels of <code>x.fac</code> must be in the order in which they are to be plotted - if they are dates, then they should be in the form <code>yyyymmdd</code> which can be achieved using <code>as.Date</code> . |
| <code>levels.length</code> | The maximum number of characters from the levels of factors to use in the row and column labels of the tables of pairwise differences and their p-values and standard errors. |
| <code>pairwise</code> | A logical indicating whether all pairwise differences of the predictions and their standard errors and p-values are to be computed and stored. If <code>FALSE</code> , the components <code>differences</code> and <code>p.differences</code> will be <code>NULL</code> in the returned <code>alldiffs</code> object. |
| <code>alpha</code> | The significance level for an LSD to compare a pair of predictions. |
| <code>inestimable.rm</code> | A logical indicating whether rows for predictions that are not estimable are to be removed from the components of the <code>alldiffs</code> object. |

Value

An `alldiffs` object that is a list with components `predictions` containing the predictions and their standard errors, `differences` containing all pairwise differences between the predictions, `p.differences` containing p-values for all pairwise differences between the predictions, `sed` containing the standard errors of all pairwise differences between the predictions, and, an `LSD` containing the mean, minimum and maximum LSDs.

See Also

`asremlPlus-package`, `alldiffs`, `print.alldiffs`, `predictionplot.asreml`, `predictparallel.asreml`, `pred.present.asreml`

Examples

```
## Not run:
Var.pred <- predict(current.asr, classify="Variety", sed=TRUE)$predictions
wald.tab <- current.asrt$wald.tab
den.df <- wald.tab[match("Variety", rownames(wald.tab)), "denDF"]
Var.diffs <- alldiffs(predictions = Var.pred$pvals,
                     sed = Var.pred$sed,
                     tdf = den.df)
Var.diffs <- predictiondiffs.asreml(classify = "Variety",
                                   alldiffs.obj=Var.diffs)
print.alldiffs(Var.diffs, which="differences")

## End(Not run)
```

`predictionplot.asreml` *This function plots the predictions for a term, possibly with error bars.*

Description

This function plots the predictions `y` that are based on `classify` stored in the data.frame `data`. The package `ggplot2` is used to produce the plots. Line plots are produced when variables involving `x.num` or `x.fac` are involved in `classify` for the predictions; otherwise, bar charts are produced. Further, for line charts, the argument `panels` determines whether a single plot or multiple plots in a single window are produced; for bar charts, the argument `panels` is ignored.

Usage

```
predictionplot.asreml(classify, y, data,
                     x.num = NULL, x.fac = NULL, nonx.fac.order = NULL,
                     colour.scheme = "colour", panels = "multiple",
                     graphics.device = NULL,
                     error.intervals = "Confidence", titles = NULL,
                     y.title = NULL, filestem = NULL, ...)
```

Arguments

| | |
|-----------------------|--|
| <code>classify</code> | a character string giving the combinations of the independent variables on which the predictions are based. It is an interaction type term formed from the independent variables, that is, separating the variable names with the <code>:</code> operator. |
|-----------------------|--|

| | |
|------------------------------|--|
| <code>y</code> | a character string giving the name of the variable that is to be plotted on the Y axis. |
| <code>data</code> | a <code>data.frame</code> containing the values of the variables to be plotted. It should be consistent with an object of class <code>asremlPredict</code> such as is stored in the <code>pvals</code> component of the <code>predictions</code> component of the value produced by <code>predict.asreml</code> ; that is, in addition to variables classifying the predictions, it will include a column with the name specified in the <code>y</code> argument, usually <code>predicted.value</code> or <code>backtransformed.predictions</code> ; each row contains a single predicted value. If <code>error.intervals</code> is not "none", then the <code>predictions</code> component and, if present, the <code>backtransforms</code> component should contain columns for the lower and upper values of the limits for the interval with names that begin with <code>lower</code> and <code>upper</code> , respectively. The second part of the name must be one of <code>Confidence</code> , <code>StandardError</code> or <code>halfLeastSignificant</code> . The last part needs to be consistent between the lower and upper limits. |
| <code>x.num</code> | A character string giving the name of the numeric covariate that corresponds to <code>x.fac</code> , is potentially included in terms in the fitted model and which corresponds to the x-axis variable. It should have the same number of unique values as the number of levels in <code>x.fac</code> . |
| <code>x.fac</code> | A character string giving the name of the factor that corresponds to <code>x.num</code> , is potentially included in terms in the fitted model and which corresponds to the x-axis variable. It should have the same number of levels as the number of unique values in <code>x.num</code> . The levels of <code>x.fac</code> must be in the order in which they are to be plotted - if they are dates, then they should be in the form <code>yyyymmdd</code> which can be achieved using <code>as.Date</code> . |
| <code>nonx.fac.order</code> | A character vector giving the order in which factors other than <code>x.fac</code> are to be plotted in faceted plots (i.e. where the number of non x factors is greater than 1). The first factor in the vector will be plotted on the X axis (if there is no <code>x.num</code> or <code>x.fac</code> . Otherwise, the order of plotting the factors is in columns (X facets) and then rows (Y facets). By default the order is in decreasing order for the numbers of levels of the non x factors. |
| <code>colour.scheme</code> | A character string specifying the colour scheme for the plots. The default is "colour" which produces coloured lines and bars, a grey background and white gridlines. A value of "black" results in black lines, grey bars and gridlines and a white background. |
| <code>panels</code> | Possible values are "single" and "multiple". When line plots are to be produced, because variables involving <code>x.num</code> or <code>x.fac</code> are involved in classify for the predictions, <code>panels</code> determines whether or not a single panel or multiple panels in a single window are produced. The <code>panels</code> argument is ignored for bar charts. |
| <code>graphics.device</code> | A character specifying a graphics device for plotting. The default is <code>graphics.device = NULL</code> , which will result in plots being produced on the current graphics device. Setting it to "windows", for example, will result in a windows graphics device being opened. |
| <code>error.intervals</code> | A character string indicating the type of error interval, if any, to calculate and plot in order to indicate uncertainty in the results. Possible values are "none", "StandardError", "Confidence" and "halfLeastSignificant". Here, any option other than "none" will result in the interval limits contained in <code>data</code> being plotted. |

| | |
|----------|--|
| titles | A list, each component of which is named for an object name and contains a character string giving a title to use in output (e.g. tables and graphs) for the object. Here they will be used for axis labels. |
| filestem | A character sting giving the beginning of the name of the file in which to save the plot. If <code>filestem = NULL</code> , the plot is not saved. The remainder of the file name will be generated automatically and consists of the following elements separated by full stops: the classify term, Bar or Line and, if <code>error.intervals</code> is not "none", one of SE, CI or LSI. The file will be saved as a 'png' file in the current work directory. |
| y.title | The title to be displayed on the y axis of any plot. |
| ... | further arguments passed to <code>ggplot</code> . |

Value

no values are returned.

See Also

[predictiondiffs.asreml](#), [pred.present.asreml](#), [ggplot](#), [Devices](#)

Examples

```
## Not run:
current.asr <- asreml(fixed = log.Turbidity ~ Benches + Sources + Type + Species +
                     Sources:Type + Sources:Species +
                     Sources:xDay + Species:xDay + Species:Date,
                     data = Runoff.longi.dat, keep.order = TRUE)
predictions <- predict(current.asr, class="Species:Date:xDay",
                      present=c("Type", "Species", "Sources"),
                      levels=list(xDay=unique(Runoff.longi.dat$xDay)))$predictions$pvals
predictions <- predictions[predictions$est.status == "Estimable",]
predictionplot.asreml(classify="Species:Date:xDay", diffs$predictions, current.asrt$wald.tab,
                     x.num = "xDay", x.fac = "Date",
                     x.title = "Days since first observation",
                     y.title = "Predicted log(Turbidity)",
                     present = c("Type", "Species", "Sources"))

diffs <- predictparallel.asreml(classify="Species:Date:xDay",
                               present=c("Type", "Species", "Sources"),
                               asreml.obj = current.asr,
                               x.num = "xDay", x.fac = "Date",
                               x.pred.values=sort(unique(Runoff.longi.dat$xDay)),
                               x.plot.values=c(0,28,56,84))
predictionplot.asreml(diffs$predictions, current.asrt$wald.tab,
                     x.num = "xDay", x.fac = "Date",
                     x.title = "Days since first observation",
                     y.title = "Predicted log(Turbidity)")

## End(Not run)
```

`predictparallel.asreml`

Uses an asreml object and a wald.tab to form the predictions and associated statistics for a term. It stores the results in an object of class alldiffs and may print the results. It can be used when a numeric vector and a factor that has parallel values both occur in the model and need to be taken into account.

Description

This function forms the predictions for term using `classify` and the supplied `asreml` object and stores them in an `alldiffs` object. If `x.num` is supplied, the predictions will be obtained for the values supplied in `x.pred.values` and, if supplied, `x.plot.values` will replace them in the `alldiffs` object that is returned. If `x.fac`, but not `x.num`, is specified, predictions will involve it and, if supplied, `x.plot.values` will replace the levels of `x.fac` in the `alldiffs` object that is returned. In order to get the correct predictions you may need to supply additional arguments to predict through ... e.g. `present`. Any aliased predictions will be removed, as will any standard error of pairwise differences involving them.

Also calculated are the approximate degrees of freedom of the standard errors of the predictions. If the denominator degrees of freedom for term are available in `wald.tab`, they are used. Otherwise the residual degrees of freedom or the maximum of the denominator degrees in `wald.tab`, excluding the Intercept, are used. Which is used depends on the setting of `ddf.na`. These degrees of freedom are used for the t-distribution on which p-values and confidence intervals are based. It is stored as an attribute to the `alldiffs` object. The degrees of freedom are also used in calculating the minimum, mean and maximum LSD for comparing pairs of predictions, which are also stored in the `alldiffs` object.

If `pairwise = TRUE`, all pairwise differences between the predictions, their standard errors, p-values and LSD statistics are computed using `predictiondiffs.asreml`. This adds them to the `alldiffs` object as additional list components named `differences`, `sed`, `p.differences` and `LSD`.

If a transformation has been applied (any one of `transform.power` is not one, `scale` is not one and `offset` is nonzero), the back-transforms of the predicted values and their lower and upper confidence intervals are added to a `data.frame` that is consistent with an object of class `asremlPredict`, such as is stored in the `pvals` component of the `prediction` component of the value produced by `predict.asreml`. This `data.frame` is added to the `alldiffs` object as a list component called `backtransforms`.

The printing of the components produced is controlled by the `tables` argument.

Usage

```
predictparallel.asreml(classify, term = NULL, asreml.obj = NULL, titles = NULL,
  x.num = NULL, x.fac = NULL,
  x.pred.values = NULL, x.plot.values = NULL,
  error.intervals = "Confidence", avsed.tolerance = 0.25,
  pairwise = TRUE, tables = "all", levels.length = NA,
  transform.power = 1, offset = 0, scale = 1,
  inestimable.rm = TRUE, wald.tab = NULL,
  alpha = 0.05, ddf.na = "residual", ddf.values = NULL,
  trace = FALSE, ...)
```

Arguments

| | |
|------------------------------|---|
| <code>classify</code> | a character string giving the variables that define the margins of the multiway table to be predicted. Multiway tables are specified by forming an interaction type term from the classifying variables, that is, separating the variable names with the <code>:</code> operator. |
| <code>term</code> | a character string giving the variables that define the term that was fitted using <code>asreml</code> and that corresponds to <code>classify</code> . It only needs to be specified when it is different to <code>classify</code> . |
| <code>asreml.obj</code> | <code>asreml</code> object for a fitted model. |
| <code>titles</code> | A list, each component of which is named for an object name and contains a character string giving a title to use in output (e.g. tables and graphs) for the object. Here they will be used for table headings. |
| <code>x.num</code> | A character string giving the name of the numeric covariate that (i) corresponds to <code>x.fac</code> , (ii) is potentially included in terms in the fitted model, and (iii) which corresponds to the x-axis variable. It should have the same number of unique values as the number of levels in <code>x.fac</code> . |
| <code>x.fac</code> | A character string giving the name of the factor that (i) corresponds to <code>x.num</code> , (ii) is potentially included in terms in the fitted model, and (iii) which corresponds to the x-axis variable. It should have the same number of levels as the number of unique values in <code>x.num</code> . The levels of <code>x.fac</code> must be in the order in which they are to be plotted - if they are dates, then they should be in the form <code>yyyymmdd</code> which can be achieved using <code>as.Date</code> . |
| <code>x.pred.values</code> | The values of <code>x.num</code> for which predicted values are required. |
| <code>x.plot.values</code> | The actual values to be plotted on the x axis. They are needed when values different to those in <code>x.num</code> are to be plotted or <code>x.fac</code> is to be plotted because there is no <code>x.num</code> term corresponding to the same term with <code>x.fac</code> . |
| <code>error.intervals</code> | A character string indicating the type of error interval, if any, to calculate and plot in order to indicate uncertainty in the results. Possible values are "none", "StandardError", "Confidence" and "halfLeastSignificant". The default is for confidence limits to be used. The "halfLeastSignificant" option results in half the mean Least Significant Difference (LSD) being added and subtracted to the predictions, the mean LSD being calculated using the average of the standard errors of all pairwise differences (SEDs) between the predictions. However, if the range of the SEDs divided by the average of the SEDs exceeds <code>aved.tolerance</code> , calculations and plotting will revert to confidence intervals. Also, half LSDs cannot be used for backtransformed values and so confidence intervals will be used instead. |
| <code>aved.tolerance</code> | The values of the range of the SEDs divided by the average of the SEDs that, if exceeded, will cause calculations and plotting to revert to confidence intervals. It should be a value between 0 and 1. |
| <code>pairwise</code> | A logical indicating whether all pairwise differences of the predictions and their standard errors and p-values are to be computed and stored. If <code>tables</code> is equal to "differences" or "all" or <code>error.intervals</code> is equal to "halfLeastSignificant", they will be stored irrespective of the value of <code>pairwise</code> . |
| <code>tables</code> | A character vector containing a combination of none, predictions, backtransforms, differences, p.differences, sed, LSD and all. These nominate which components of the <code>alldiffs</code> object to print. |

| | |
|-----------------|---|
| levels.length | The maximum number of characters from the the levels of factors to use in the row and column labels of the tables of pairwise differences and their p-values and standard errors. |
| transform.power | A number specifying the power of a transformation, if one has been applied to the response variable. Unless it is equal to 1, the default, back-transforms of the predictions will be obtained and presented in tables or graphs as appropriate. The back-transformation raises the predictions to the power equal to the reciprocal of transform.power, unless it equals 0 in which case the exponential of the predictions is taken. |
| offset | A number that has been added to each value of the response after any scaling and before applying any power transformation. |
| scale | A number by which each value of the response has been multiply before adding any offset and applying any power transformation. |
| inestimable.rm | A logical indicating whether rows for predictions that are not estimable are to be removed from the components of the <code>alldiffs</code> object. |
| wald.tab | a data frame containing the pseudo-anova table for the fixed terms produced by a call to <code>wald.asreml</code> . The main use of it here is in determinining the degrees of freedom of the standard errors of the predictions. denominator degrees of freedom when p-values or confidence intervals are to be calculated. |
| alpha | the significance level for a test or one minus the confidence level for confidence intervals. |
| dDF.na | the method to use to obtain approximate denominator degrees of freedom. when the numeric or algebraic methods produce an NA. Consistent with when no denDF are available, the default is "residual" and so the residual degrees of freedom from <code>asreml.obj\$nedf</code> are used. If <code>dDF.na = "none"</code> , no substitute denominator degrees of freedom are employed; if <code>dDF.na = "maximum"</code> , the maximum of those denDF that are available, excluding that for the Intercept, is used; if all denDF are NA, <code>asreml.obj\$nedf</code> is used. If <code>dDF.na = "supplied"</code> , a vector of values for the denominator degrees of freedom is to be supplied in <code>dDF.values</code> . Any other setting is ignored and a warning message produced. Generally, substituting these degrees of freedom is anticonservative in that it is likely that the degrees of freedom used will be too large. |
| dDF.values | A vector of values to be used when <code>dDF.na = "supplied"</code> . Its values will be used when denDF in a test for a fixed effect is NA. This vector must be the same length as the number of fixed terms, including (Intercept) whose value could be NA. |
| trace | if TRUE then partial iteration details are displayed when ASReml-R functions are invoked; if FALSE then no output is displayed. |
| ... | further arguments passed to <code>predict.asreml</code> . |

Value

An `alldiffs` object with predictions and their standard errors and, depending on the settings of the arguments, all pairwise differences between predictions, their standard errors and p-values and LSD statistics. If `power.transform` is not one, it will contain a data.frame with the backtransformed predictions. If `error.intervals` is not "none", then the predictions component and, if present, the backtransforms component will contain columns for the lower and upper values of the limits for the interval. The names of these columns will consist of three parts separated by full stops: 1)

the first part will be lower or upper; 2) the second part will be one of Confidence, StandardError or halfLeastSignificant; 3) the third component will be limits.

The name of the response, the term, the classify and tdf, as well as the degrees of freedom of the standard error, will be set as attributes to the object.

See Also

[alldiffs](#), [print.alldiffs](#), [predictiondiffs.asreml](#), [pred.present.asreml](#), [as.Date](#), [predictionplot.asreml](#), [pred.present.asreml](#), [predict.asreml](#)

Examples

```
## Not run:
current.asr <- asreml(fixed = pH ~ Benches + (Sources * (Type + Species)),
                     random = ~ Benches:MainPlots,
                     keep.order=TRUE, data= Runoff.longi.dat)
diffs <- predictparallel.asreml(classify = "Sources:Type",
                               asreml.obj = current.asr,
                               x.num = "xDay", x.fac = "Date",
                               x.pred.values=sort(unique(Runoff.longi.dat$xDay)),
                               x.plot.values=c(0,28,56,84))

## End(Not run)
```

print.alldiffs

Prints the values in an [alldiffs](#) object in a nice format.

Description

Prints the predictions and standard errors from using asreml to fit models in the same way as asreml prints them. Also prints out all pairwise differences between the predictions to 2 significant figures, along with their p-values and standard errors to 4 decimal places. If LSDs are requested the mean, minimum and maximum LSDs will be printed.

Usage

```
## S3 method for class 'alldiffs'
print(x, which = "all", ...)
```

Arguments

| | |
|-------|--|
| x | A list returned from asreml.predictiondiffs. |
| which | A character vector containing a combination of predictions, backtransforms, differences, p.differences, sed, LSD and all. These nominate which components of the alldiffs object to print. |
| ... | further arguments passed to or from other methods. |

Value

No value is returned, but the elements of the list in x, returned from asreml.predictiondiffs, are printed.

See Also

[alldiffs](#), [predictiondiffs.asreml](#)

Examples

```
## Not run:
print.alldiffs(predictiondiffs.asreml(asreml=splitplot.asr, wald=wald,
                                     present=c("Type", "Sources", "Species"),
                                     factors=list("Species")))

## End(Not run)
```

| | |
|----------------|--|
| print.asrtests | <i>Prints the values in an asrtests object</i> |
|----------------|--|

Description

Prints a summary of the asreml object and the test.summary data.frame that are stored in the [asrtests](#) object.

Usage

```
## S3 method for class 'asrtests'
print(x, which = "all", ...)
```

Arguments

| | |
|-------|--|
| x | An asrtests object. |
| which | Which elements of the asrtests object to print. Possible values are some combination of asremlsummary, pseudoanova, testsummary and all. |
| ... | further arguments passed to print. |

Value

No value is returned, but the elements of the list in x are printed.

See Also

[asrtests](#), [asremlPlus-package](#)

Examples

```
## Not run:
data(Wheat.dat)
current.asr <- asreml(yield ~ Rep + WithinColPairs + Variety,
                    random = ~ Row + Column + units,
                    rcov = ~ ar1(Row):ar1(Column),
                    data=Wheat.dat)
current.asrt <- asrtests(current.asr, NULL, NULL)
current.asrt <- rmboundary.asrtests(current.asrt)
# Test Row autocorrelation
current.asrt <- testrcov.asrtests("~ Row:ar1(Column)", current.asrt,
```

```

                                label="Row autocorrelation", simpler=TRUE)
print(current.asrt)

## End(Not run)

```

```
recalc.wald.tab.asrtests
```

Recalculates the denDF, F.inc and P values for a table of Wald test statistics obtained using wald.asrem1

Description

If some or all denDF are not available, either because they are NA or because F.inc values were not calculated, this function allows the user to specify how a approximate denDF values are to be obtained. This is done through the dDF.na and dDF.values arguments. Note that if denDF values are available in the Wald table then only those that are NA will be replaced.

Usage

```

recalc.wald.tab.asrtests(asrtests.obj, recalc.wald = FALSE,
                        denDF="default", dDF.na = "none",
                        dDF.values = NULL, trace = FALSE, ...)

```

Arguments

- | | |
|--------------|--|
| asrtests.obj | an asrtests object for a fitted model that is a list containing an asrem1 object, a wald.tab data.frame with 4 columns, and a data.frame with 5 columns that records any previous changes and tests in the fitted model. |
| recalc.wald | a logical indicating whether to call wald.asrem1 to recalculate the pseudoanova table for the model fit stored in the asrem1 object contained in asrtests . |
| denDF | Specifies the method to use in computing approximate denominator degrees of freedom when wald.asrem1 is called. Can be none to suppress the computations, numeric for numerical methods, algebraic for algebraic methods or default, the default, to automatically choose numeric or algebraic computations depending on problem size. The denominator degrees of freedom are calculated according to Kenward and Roger (1997) for fixed terms in the dense part of the model. |
| dDF.na | the method to use to obtain substitute denominator degrees of freedom. when the numeric or algebraic methods produce an NA. If dDF.na = "none", no substitute denominator degrees of freedom are employed; if dDF.na = "residual", the residual degrees of freedom from asrem1.obj\$nedf are used; if dDF.na = "maximum", the maximum of those denDF that are available, excluding that for the Intercept, is used; if all denDF are NA, asrem1.obj\$nedf is used. If dDF.na = "supplied", a vector of values for the denominator degrees of freedom is to be supplied in dDF.values. Any other setting is ignored and a warning message produced. Generally, substituting these degrees of freedom is anticonservative in that it is likely that the degrees of freedom used will be too large. |
| dDF.values | A vector of values to be used when dDF.na = "supplied". Its values will be used when denDF in a test for a fixed effect is NA. This vector must be the same length as the number of fixed terms, including (Intercept) whose value could be NA. |

`trace` if TRUE then partial iteration details are displayed when ASReml-R functions are invoked; if FALSE then no output is displayed.

`...` further arguments passed to `asreml` and to `wald.asreml`.

Value

A `wald.tab`: a 4-column data.frame containing a pseudo-anova table for the fixed terms produced by `wald.asreml`.

See Also

[asrtests](#), [testranfix.asrtests](#)

Examples

```
## Not run:
wald.tab <- recalc.wald.tab.asrtests(current.asrt,
                                     dDF.na = "supplied",
                                     dDF.values = c(NA,rep(c(330,346), c(4,3))))

## End(Not run)
```

| | |
|------------------------------|---|
| <code>reml.lrt.asreml</code> | <i>Performs REML likelihood ratio test.</i> |
|------------------------------|---|

Description

Extracts the REML log likelihood and number of variance parameters from two `asreml` objects. It assumes that the second `asreml` object is the result of fitting a model that is a reduced version of the model for the first object. In the case where the reduced model is obtained by setting positively-constrained variance parameters in the full model to zero, the `positive.zero` argument should be set to TRUE so that the p-value is computed using a mixture of chi-square distributions as described in Self and Liang (1987).

The function checks that the models do not differ in either their fixed or sparse models.

The function `reml.lrt` is provided for backwards compatibility.

Usage

```
reml.lrt.asreml(full.asreml.obj, reduced.asreml.obj, positive.zero=FALSE)
```

Arguments

`full.asreml.obj` `asreml` object for the full model.

`reduced.asreml.obj` `asreml` object for the reduced model.

`positive.zero` Indicates whether the hypothesized values for the variance components being tested are on the boundary of the parameter space. For example, this is true for positively-constrained variance components that, under the reduced model, are zero.

Value

A data frame containing the log of the likelihood ratio, its degrees of freedom and its p-value.

Note

The degrees of freedom for the test are computed as the difference between the two models in the number of variance parameters that are unfixed, nonsingular and unconstrained by relationships among them.

This procedure is only appropriate when the null hypothesis is that all parameters are on the boundary of the parameter space or that all parameters are in the interior of the parameter space. Mixed cases have been discussed by Self and Liang (1987), but are not implemented here.

References

Self, S.G., and Liang, K-Y. (1987) Asymptotic Properties of Maximum Likelihood Estimators and Likelihood Ratio Tests Under Nonstandard Conditions. *Journal of the American Statistical Association*, **82**, 605-10.

See Also

[info.crit.asreml](#), [testranfix.asrtests](#)

Examples

```
## Not run:
  reml.lrt.asreml(ICV.max, ICV.red, positive.zero=TRUE)

## End(Not run)
```

| | |
|---------------------|--|
| rmboundary.asrtests | <i>Removes any boundary or singular variance components from the fit stored in asreml.obj and records their removal in a data.frame.</i> |
|---------------------|--|

Description

Any terms specified in the random model that are estimated on the boundary or are singular and can be removed are removed from the fit stored in an asreml object. Terms that specify multiple parameters in the random model cannot be removed (e.g. terms specified using the at function with more than one level of the factor) and terms in rcov model are not removed. Terms that can be removed are selected for removal in the following order based on whether they involve: (i) a dev function, (ii) only factors, (iii) an spl function, (iv) a pol function and (v) a lin function or a variable that is an integer or a numeric. It should be noted that this order of removal presumes that random deviation terms are specified via the dev function rather than via a random factor. Once the earliest of these above classes with a boundary term is identified, a term within this class is selected for removal. For all classes, except for factor-only terms, the smallest term with the largest number of variables/factors is removed. Amongst factor-only terms, the smallest term with the smallest number of variables/factors is removed. After each variance component is removed, a row for it is added to the test.summary data.frame and the model refitted. If there are further boundary or singular terms, one is removed using the above strategy. This process continues until there are no further boundary or singular variance components that are removable. Other types of boundary or singular terms, which cannot be removed, are reported in warning messages.

Usage

```
rmboundary.asrtests(asrtests.obj, trace = FALSE, update = TRUE,
  set.terms = NULL, ignore.suffices = TRUE,
  constraints = "P", initial.values = NA, ...)
```

Arguments

| | |
|------------------------------|---|
| <code>asrtests.obj</code> | an asrtests object for a fitted model that is a list containing an <code>asreml</code> object, a <code>wald.tab</code> data.frame with 4 columns, and a data.frame with 5 columns that records any previous changes and tests in the fitted model. |
| <code>trace</code> | if TRUE then partial iteration details are displayed when ASReml-R functions are invoked; if FALSE then no output is displayed. |
| <code>update</code> | if TRUE then <code>update.asreml</code> is called to fit the model with any boundary terms removed. In doing this the arguments <code>R.param</code> and <code>G.param</code> are set to those in the <code>asreml</code> object stored in <code>asrtests.obj</code> so that the values from the previous model are used as starting values. If FALSE then a call is made to <code>asreml</code> in which the only changes from the previous call are that (i) the terms for boundary variance components are removed from the models and (ii) modifications specified via ... are made. |
| <code>set.terms</code> | a character vector specifying the terms that are to have constraints and/or initial values set prior to fitting. |
| <code>ignore.suffices</code> | a logical vector specifying whether the suffices of the <code>asreml</code> -assigned names of the variance terms (i.e. the information to the right of an "!", other than "R!") is to be ignored in matching elements of terms. If TRUE for an element of terms, the suffices are stripped from the <code>asreml</code> -assigned names. If FALSE for an element of terms, the element must exactly match an <code>asreml</code> -assigned name for a variance term. This vector must be of length one or the same length as <code>terms</code> . If it is of length one then the same action is applied to the <code>asreml</code> -assigned suffices for all the terms in <code>terms</code> . |
| <code>constraints</code> | a character vector specifying the constraints to be applied to the terms specified in <code>terms</code> . This vector must be of length one or the same length as <code>terms</code> . If it is of length one then the same constraint is applied to all the terms in <code>terms</code> . If any of the constraints are equal to NA then they are left unchanged for those terms. |
| <code>initial.values</code> | a character vector specifying the initial values for the terms specified in <code>terms</code> . This vector must be of length one or the same length as <code>terms</code> . If it is of length one then the same initial value is applied to all the terms in <code>terms</code> . If any of the <code>initial.values</code> are equal to NA then they are left unchanged for those terms. |
| <code>...</code> | further arguments passed to <code>asreml</code> . |

Value

An [asrtests](#) object, which is a list containing:

1. `asreml.obj`: an `asreml` object containing the fit of the model after all boundary and singular terms have been removed;
2. `wald.tab`: a 4-column data.frame containing a pseudo-anova table for the fixed terms produced by `wald.asreml`;
3. `test.summary`: a data.frame with columns `term`, `DF`, `denDF`, `p` and `action`. A row is added to it for each of the boundary terms removed, the row containing the name of the term, one for the DF, NA for the p-value and Boundary for the action.

See Also

[asrtests](#), [addrm.terms.asrtests](#), [testranfix.asrtests](#), [testrcov.asrtests](#),
[newfit.asreml](#), [sig.devn.reparam.asrtests](#), [choose.model.asrtests](#)

Examples

```
## Not run:
current.asrt <- rmboundary.asrtests(current.asrt)

## End(Not run)
```

```
setvarianceterms.asreml
```

allows the setting of constraints and initial values for terms in the random and rcov arguments of an asreml call, with the resulting call being evaluated.

Description

Takes an unevaluated call and evaluates the call after setting the constraints and initial values for the terms specified in terms. The elements of terms are matched with those generated by asreml and used, for example, in the varcomp component of a summary.asreml object. These names generally include descriptive suffices. To match an element of terms that includes such a suffix, set ignore.suffices to FALSE so that a literal match between the element and the assigned names is sought.

Usage

```
setvarianceterms.asreml(call, terms, ignore.suffices = TRUE,
                        constraints = "P", initial.values = NA, ...)
```

Arguments

| | |
|-----------------|---|
| call | an unevaluated call to asreml. One way to create such a call is to use the call function with its name argument set to "asreml". Another is to obtain it from the call component of an asreml object (e.g. call <- asreml.obj\$call). |
| terms | a character vector specifying the terms that are to have constraints and/or initial values specified. |
| ignore.suffices | a logical vector specifying whether the suffices of the asreml-assigned names of the variance terms (i.e. the information to the right of an "!", other than "R!") is to be ignored in matching elements of terms. If TRUE for an element of terms, the suffices are stripped from the asreml-assigned names. If FALSE for an element of terms, the element must exactly match an asreml-assigned name for a variance term. This vector must be of length one or the same length as terms. If it is of length one then the same action is applied to the asreml-assigned suffices for all the terms in terms. |
| constraints | a character vector specifying the constraints to be applied to the terms specified in terms. This vector must be of length one or the same length as terms. If it is of length one then the same constraint is applied to all the terms in terms. If any of the constraints are equal to NA then they are left unchanged for those terms. |

`initial.values` a character vector specifying the initial values for the terms specified in `terms`. This vector must be of length one or the same length as `terms`. If it is of length one then the same initial value is applied to all the terms in `terms`. If any of the `initial.values` are equal to NA then they are left unchanged for those terms.

... additional arguments to be added to the call, or arguments in the call with changed values.

Value

An `asreml` object.

References

Butler, D. G., et al. (2010). *Analysis of Mixed Models for S language environments: ASReml-R reference manual*. Brisbane, DPI Publications.

See Also

`update.asreml`

Examples

```
## Not run:
m1.call <- call("asreml",
  fixed = Height ~ (Block + Irrig)*csDay.num,
  random= ~ spl(csDay.num)/(Irrig+Block)
  + dev(csDay.num)
  + str(~Block:Plot/csDay.num, ~us(2):id(20))
  + Block:Plot:spl(csDay.num),
  data=quote(dat)) ##use quote to stop evaluation of dat here
terms <- c("Block:Plot+Block:Plot:csDay.num!us(2).2:1", "R!variance")
m1.asreml <- setvarianceterms.asreml(m1.call, terms,
  constraints=c("U", "P"),
  initial=c(NA,3),
  ignore.suffices=c(FALSE,TRUE))

summary(m1.asreml)

## End(Not run)
```

sig.devn.reparam.asrtests

This function reparameterizes each random (deviations) term involving devn.fac to a fixed term and ensures that the same term, with trend.num replacing devn.fac, is included if any other term with trend.num is included in terms.

Description

This function reparameterizes each random (deviations) term involving `devn.fac` to a fixed term and ensures that the same term with `trend.num` replacing `devn.fac` is included if any other term with `trend.num` is included in `terms`. It also ensures that any term with `spl{trend.num}` replacing `devn.fac` in a term being reparameterized is removed from the model.

Usage

```
sig.devn.reparam.asrtests(terms = NULL, asrtests.obj,
                           trend.num = NULL, devn.fac = NULL,
                           denDF = "default", trace = FALSE, update = TRUE,
                           set.terms = NULL, ignore.suffices = TRUE,
                           constraints = "P", initial.values = NA,...)
```

Arguments

| | |
|------------------------------|---|
| <code>terms</code> | a character string vector giving the terms that are to be reparameterized. |
| <code>asrtests.obj</code> | an <code>asrtests</code> object for a fitted model that is a list containing an <code>asreml</code> object, a <code>wald.tab</code> data.frame with 4 columns, and a data.frame with 5 columns that records any previous changes and tests in the fitted model. |
| <code>trend.num</code> | A character string giving the name of the numeric covariate that corresponds to <code>devn.fac</code> and is potentially included in terms in the fitted model. |
| <code>devn.fac</code> | A character string giving the name of the factor that corresponds to <code>trend.num</code> and is included in terms in the fitted model. |
| <code>denDF</code> | Specifies the enthod to use in computing approximate denominator degrees of freedom when <code>wald.asreml</code> is called. Can be none to suppress the computations, numeric for numerical methods, algebraic for algebraic methods or default, the default, to autommatically choose numeric or algebraic computations depending on problem size. The denominator degrees of freedom are calculated according to Kenward and Roger (1997) for fixed terms in the dense part of the model. |
| <code>trace</code> | if TRUE then partial iteration details are displayed when ASReml-R functions are invoked; if FALSE then no output is displayed. |
| <code>update</code> | if TRUE then <code>update.asreml</code> is called in removing and adding terms to the model. In doing this the arguments <code>R.param</code> and <code>G.param</code> are set to those in the <code>asreml</code> object stored in the supplied <code>asrtests.obj</code> so that the values from the previous model are used as starting values. If FALSE then calls are made to <code>asreml</code> in which the only changes from the previous call are (i) that the models are updated and (ii) modifications specified via . . . are made. |
| <code>set.terms</code> | a character vector specifying the terms that are to have constraints and/or initial values set prior to fitting. |
| <code>ignore.suffices</code> | a logical vector specifying whether the suffices of the <code>asreml</code> -assigned names of the variance terms (i.e. the information to the right of an "!", other than "R!") is to be ignored in matching elements of terms. If TRUE for an element of terms, the suffices are stripped from the <code>asreml</code> -assigned names. If FALSE for an element of terms, the element must exactly match an <code>asreml</code> -assigned name for a variance term. This vector must be of length one or the same length as terms. If it is of length one then the same action is applied to the <code>asreml</code> -assigned suffices for all the terms in terms. |
| <code>constraints</code> | a character vector specifying the constraints to be applied to the terms specified in terms. This vector must be of length one or the same length as terms. If it is of length one then the same constraint is applied to all the terms in terms. If any of the constraints are equal to NA then they are left unchanged for those terms. |
| <code>initial.values</code> | a character vector specifying the initial values for the terms specified in terms. This vector must be of length one or the same length as terms. If it is of length |

one then the same initial value is applied to all the terms in terms. If any of the initial.values are equal to NA then they are left unchanged for those terms.

... further arguments passed to asreml via [addrm.terms.asrtests](#).

Value

An [asrtests](#) object, which is a list containing:

1. `asreml.obj`: an `asreml` object containing the fit of the model after all boundary and singular terms have been removed;
2. `wald.tab`: a 4-column data.frame containing a pseudo-anova table for the fixed terms produced by `wald.asreml`;
3. `test.summary`: a data.frame with columns term, DF, denDF, p and action. It contains a row for each term that is dropped, added or tested or a note that several terms have been added or removed.

See Also

[asrtests](#), [addrm.terms.asrtests](#), [testranfix.asrtests](#), [testrcov.asrtests](#), [newfit.asreml](#), [choose.model.asrtests](#)

Examples

```
## Not run:
current.asrt <- sig.devn.reparam.asrtests(choose$`sig.terms`, current.asrt,
                                         trend.num = "xDay", devn.fac = "Date",
                                         denDF = "algebraic")

## End(Not run)
```

| | |
|------------------------------|--|
| <code>simulate.asreml</code> | <i>Produce sets of simulated data from a multivariate normal distribution and save quantiles related to the simulated data</i> |
|------------------------------|--|

Description

Produce a set of simulated data corresponding to an `asreml` model, along with its fitted values and residuals. A variance matrix `V`, corresponding to the random and `rcov` models must be supplied. What to save is specified by the `which` argument.

Usage

```
## S3 method for class 'asreml'
simulate(object, nsim=100, seed = NULL, means=NULL, V, tolerance = 1E-10,
        update = TRUE, trace = FALSE, which="data", units = "ignore", ...)
```

Arguments

| | |
|-----------|---|
| object | An asreml object from a call to asreml in which the data argument has been set. |
| means | The vector of means to be used in generating simulated data sets. If it is NULL, the fitted values based on object are used. |
| V | The fitted variance matrix, i.e. having the pattern and values that conform to the model fit stored in the supplied object. |
| nsim | The number of data sets to be simulated. |
| seed | A single value, interpreted as an integer, that specifies the starting value of the random number generator. |
| tolerance | The value such that eigenvalues less than it are considered to be zero. |
| update | if TRUE then the arguments R.param and G.param are set to those in the asreml object supplied in object so that the values from the original model are used as starting values. If FALSE then calls are made to asreml in which the only changes from the previous call are (i) the model is fitted to simulated data and (ii) modifications specified via ... are made, except that changes cannot be made to any of the models. |
| trace | if TRUE then partial iteration details are displayed when ASReml-R functions are invoked; if FALSE then no output is displayed. |
| which | The quantities from the simulated data set to be stored. Any combination of "response", "residuals" and "fitted", or "all". If residuals and/or fitted is specified, those for the analysis stored in object will be added to the data.frame nominated in the data argument of object and the modified data.frame added as a component named data in the list that is the value returned by the function. |
| units | A character indicating whether the BLUPs for units are added to the residuals when this reserved factor is included in the random model. Possible values are addtoresiduals and ignore. |
| ... | Other arguments that are passed down to the function asreml. Changes to the models are not allowed. Other changes are dangerous and generally should be avoided. |

Details

Generate nsim set of data and analyse them using asreml using the model in object. Note, if the analysis for a data set does not converge in maxiter iterations, it is discarded and a replacement data set generated. The value of maxiter can be specified in the call to simulate.asreml. The fitted values and residuals are extracted as required. If aom = TRUE when the simulated data are analysed, standardised conditional residuals are stored. If which includes residuals or fitted, the specified quantities for the observed data are added to the data.frame on which the fit in object is based.

Value

An list with the following components whose presence depends on the setting of which:

1. **observed:** present if which includes residuals or fitted, in which case it will be the data.frame on which the fit in object is based, with residuals and/or fitted.
2. **data:** present if which includes data, a data.frame containing the simulated data sets.

3. **fitted:** present if which includes fitted, a data.frame containing the fitted values from the analyses of the simulated data sets.
4. **residuals:** present if which includes residuals, a data.frame containing the residuals from the analyses of the simulated data sets.

Author(s)

Chris Brien

See Also

asreml, [variofaces.asreml](#), [plotvariofaces.asreml](#).

Examples

```
## Not run:
data(Wheat.dat)
current.asr <- asreml(yield ~ Rep + WithinColPairs + Variety,
                     random = ~ Row + Column + units,
                     rcov = ~ ar1(Row):ar1(Column),
                     data=Wheat.dat)
current.asrt <- asrtests(current.asr, NULL, NULL)
current.asrt <- rmboundary.asrtests(current.asrt)
# Form variance matrix based on estimated variance parameters
s2 <- current.asr$sigma2
gamma.Row <- current.asr$gammas[1]
gamma.unit <- current.asr$gammas[2]
rho.r <- current.asr$gammas[4]
rho.c <- current.asr$gammas[5]
row.ar1 <- mat.ar1(order=10, rho=rho.r)
col.ar1 <- mat.ar1(order=15, rho=rho.c)
V <- gamma.Row * fac.sumop(Wheat.dat$Row) +
    gamma.unit * diag(1, nrow=150, ncol=150) +
    mat.dirprod(col.ar1, row.ar1)
V <- s2*V

#Produce residuals from 100 simulated data sets
resid <- simulate(current.asr, V=V, which="residuals")

## End(Not run)
```

| | |
|---------------------|---|
| testranfix.asrtests | <i>Tests for a single fixed or random term in model fitted using asreml</i> |
|---------------------|---|

Description

Tests for a single term, using a REML LRT for a random term or based on Wald statistics for a fixed term. The term must be in the fitted model. A random term is removed from the model fit and a REML likelihood ratio test is performed using [reml.lrt.asreml](#). It compares the fit of the model in `asreml.obj` and the newly fitted model without the term. If the newly fitted model is retained, any boundary terms are then removed using [rmboundary.asrtests](#). For a fixed term, the probability of the Wald statistics is extracted from the pseudo-anova table produced by `wald.asreml`. If this is available in the [asrtests](#) object, it is used; otherwise `wald.asreml` is called to add it to

the `asrtests` object. Whether nonsignificant terms are dropped is controlled by `drop.ran.ns` for random terms and `drop.fix.ns` for fixed terms. A row is added to the `test.summary.data.frame` for the term that is tested.

Usage

```
testranfix.asrtests(term=NULL, asrtests.obj, alpha = 0.05,
  drop.ran.ns = TRUE, positive.zero = FALSE,
  drop.fix.ns = FALSE, denDF="default", dDF.na = "none",
  dDF.values = NULL, trace = FALSE, update = TRUE,
  set.terms = NULL, ignore.suffices = TRUE,
  constraints = "P", initial.values = NA, ...)
```

Arguments

| | |
|----------------------------|--|
| <code>term</code> | a single model term that is valid in <code>asreml</code> , stored as a character. |
| <code>asrtests.obj</code> | an <code>asrtests</code> object for a fitted model that is a list containing an <code>asreml</code> object, a <code>wald.tab</code> data.frame with 4 columns, and a data.frame with 5 columns that records any previous changes and tests in the fitted model. |
| <code>alpha</code> | the significance level for the test. |
| <code>drop.ran.ns</code> | a logical indicating whether to drop a random term from the model when it is nonsignificant. |
| <code>positive.zero</code> | a logical indicating whether the hypothesized values for the variance components being tested are on the boundary of the parameter space. For example, this is be true for positively-constrained variance components that, under the reduced model, are zero. |
| <code>drop.fix.ns</code> | a logical indicating whether to drop a fixed term from the model when it is nonsignificant |
| <code>denDF</code> | Specifies the enthod to use in computing approximate denominator degrees of freedom when <code>wald.asreml</code> is called. Can be <code>none</code> to suppress the computations, <code>numeric</code> for numerical methods, <code>algebraic</code> for algebraic methods or <code>default</code> , the default, to automatically choose numeric or algebraic computations depending on problem size. The denominator degrees of freedom are calculated according to Kenward and Roger (1997) for fixed terms in the dense part of the model. |
| <code>dDF.na</code> | the method to use to obtain substitute denominator degrees of freedom. when the numeric or algebraic methods produce an NA. If <code>dDF.na = "none"</code> , no substitute denominator degrees of freedom are employed; if <code>dDF.na = "residual"</code> , the residual degrees of freedom from <code>asreml.obj\$nedf</code> are used; if <code>dDF.na = "maximum"</code> , the maximum of those <code>denDF</code> that are available, excluding that for the Intercept, is used; if all <code>denDF</code> are NA, <code>asreml.obj\$nedf</code> is used. If <code>dDF.na = "supplied"</code> , a vector of values for the denominator degrees of freedom is to be supplied in <code>dDF.values</code> . Any other setting is ignored and a warning message produced. Generally, substituting these degrees of freedom is anticonservative in that it is likely that the degrees of freedom used will be too large. |
| <code>dDF.values</code> | A vector of values to be used when <code>dDF.na = "supplied"</code> . Its values will be used when <code>denDF</code> in a test for a fixed effect is NA. This vector must be the same length as the number of fixed terms, including (Intercept) whose value could be NA. |
| <code>trace</code> | if TRUE then partial iteration details are displayed when <code>ASReml-R</code> functions are invoked; if FALSE then no output is displayed. |

| | |
|------------------------------|---|
| <code>update</code> | if TRUE then <code>update.asreml</code> is called to fit the model to be tested. In doing this the arguments <code>R.param</code> and <code>G.param</code> are set to those in the <code>asreml</code> object stored in <code>asrtests.obj</code> so that the values from the previous model are used as starting values. If FALSE then a call is made to <code>asreml</code> in which the only changes from the previous call are that (i) models are modified for the supplied terms and (ii) modifications specified via <code>...</code> are made. |
| <code>set.terms</code> | a character vector specifying the terms that are to have constraints and/or initial values set prior to fitting. |
| <code>ignore.suffices</code> | a logical vector specifying whether the suffices of the <code>asreml</code> -assigned names of the variance terms (i.e. the information to the right of an "!", other than "R!") is to be ignored in matching elements of terms. If TRUE for an element of terms, the suffices are stripped from the <code>asreml</code> -assigned names. If FALSE for an element of terms, the element must exactly match an <code>asreml</code> -assigned name for a variance term. This vector must be of length one or the same length as <code>terms</code> . If it is of length one then the same action is applied to the <code>asreml</code> -assigned suffices for all the terms in <code>terms</code> . |
| <code>constraints</code> | a character vector specifying the constraints to be applied to the terms specified in <code>terms</code> . This vector must be of length one or the same length as <code>terms</code> . If it is of length one then the same constraint is applied to all the terms in <code>terms</code> . If any of the constraints are equal to NA then they are left unchanged for those terms. |
| <code>initial.values</code> | a character vector specifying the initial values for the terms specified in <code>terms</code> . This vector must be of length one or the same length as <code>terms</code> . If it is of length one then the same initial value is applied to all the terms in <code>terms</code> . If any of the <code>initial.values</code> are equal to NA then they are left unchanged for those terms. |
| <code>...</code> | further arguments passed to <code>asreml</code> and to <code>wald.asreml</code> . |

Value

An `asrtests` object, which is a list containing:

1. `asreml.obj`: an `asreml` object containing the fit after the term has been tested; it will be a new model if the term is nonsignificant and the appropriate argument out of `drop.ran.ns` and `drop.fix.ns` is TRUE;
2. `wald.tab`: a 4-column data.frame containing a pseudo-anova table for the fixed terms produced by `wald.asreml`;
3. `test.summary`: a data.frame with columns `term`, `DF`, `denDF`, `p` and `action`. A row is added to it for each term that is tested, the row containing the name of the term, the degrees of freedom (numerator DF for a Wald test and the number of extra parameters for a REML ratio tests), the p-value and a for the action taken. Possible codes are: Dropped, Retained, Significant, Nonsignificant, Absent, Added, Removed and Boundary. If the changed model did not converge, Unconverged will be added to the code. Note that the logical `asreml.obj$converge` also reflects whether there is convergence.

If the term is not in the model, then the supplied `asreml` object will be returned. Also, `reml.test` will have the likelihood ratio and the p-value set to NA and the degrees of freedom to zero. Similarly, the row of `test.summary` for the term will have its name, `DF` set to NA, p-value set to NA, and `action` set to Absent.

See Also

[asremlPlus-package](#), [asrtests](#), [choose.model.asrtests](#), [reml.lrt.asreml](#), [rmboundary.asrtests](#), [newfit.asreml](#), [sig.devn.reparam.asrtests](#), [addrm.terms.asrtests](#)

Examples

```
## Not run:
data(Wheat.dat)
current.asr <- asreml(yield ~ Rep + WithinColPairs + Variety,
                     random = ~ Row + Column + units,
                     rcov = ~ ar1(Row):ar1(Column),
                     data=Wheat.dat)
current.asrt <- asrtests(current.asr, NULL, NULL)
current.asrt <- rmboundary.asrtests(current.asrt)
# Test nugget term
current.asrt <- testranfix.asrtests("units", current.asrt, positive=TRUE)

## End(Not run)
```

| | |
|-------------------|--|
| testrcov.asrtests | <i>Fits a new rcov formula using asreml and tests whether the change is significant.</i> |
|-------------------|--|

Description

Fits a new rcov formula using `asreml` and tests whether the change is significant. If `simpler = FALSE` the model to be fitted must be more complex than the one whose fit has been stored in `asrtests.obj`. That is, the new model must have more parameters. However, if `simpler = TRUE` the model to be fitted must be simpler than the one whose fit has been stored in `asrtests.obj` in that it must have fewer parameters. Any boundary terms are removed using `rmboundary.asrtests`, which may mean that the models are not nested. The test is a REML likelihood ratio test that is performed using `reml.lrt.asreml`, which is only valid if the models are nested. It compares the newly fitted model with the fit of the model in `asrtest.obj`. A row is added to the `test.summary` data.frame using the supplied label.

Usage

```
testrcov.asrtests(terms=NULL, asrtests.obj, label = "R model",
                  simpler = FALSE, alpha = 0.05,
                  positive.zero = FALSE, denDF="default",
                  update = TRUE, trace = FALSE,
                  set.terms = NULL, ignore.suffices = TRUE,
                  constraints = "P", initial.values = NA, ...)
```

Arguments

| | |
|---------------------------|---|
| <code>terms</code> | a model for the <code>rcov</code> argument in <code>asreml</code> , stored as a character. |
| <code>asrtests.obj</code> | an <code>asrtests</code> object for a fitted model that is a list containing an <code>asreml</code> object, a <code>wald.tab</code> data.frame with 4 columns, and a data.frame with 5 columns that records any previous changes and tests in the fitted model. |
| <code>label</code> | a character string to use as the label in <code>test.summary</code> and which indicates what is being tested. |
| <code>simpler</code> | a logical indicating whether the new model to be fitted is simpler than the already fitted model whose fit is stored in <code>asrtests.obj</code> . |
| <code>alpha</code> | the significance level for the test. |

| | |
|-----------------|---|
| positive.zero | a logical indicating whether the hypothesized values for the variance components being tested are on the boundary of the parameter space. For example, this is be true for positively-constrained variance components that, under the reduced model, are zero. |
| denDF | Specifies the method to use in computing approximate denominator degrees of freedom when <code>wald.asreml</code> is called. Can be <code>none</code> to suppress the computations, <code>numeric</code> for numerical methods, <code>algebraic</code> for algebraic methods or <code>default</code> , the default, to automatically choose numeric or algebraic computations depending on problem size. The denominator degrees of freedom are calculated according to Kenward and Roger (1997) for fixed terms in the dense part of the model. |
| update | if <code>TRUE</code> then <code>update.asreml</code> is called to fit the model with the <code>rcov</code> model supplied in <code>terms</code> . In doing this the arguments <code>R.param</code> and <code>G.param</code> are set to those in the <code>asreml</code> object stored in <code>asrtests.obj</code> so that the values from the previous model are used as starting values. If <code>FALSE</code> then a call is made to <code>asreml</code> in which the only changes from the previous call are that (i) <code>rcov</code> model is that specified in <code>terms</code> and (ii) modifications specified via <code>...</code> are made. |
| trace | if <code>TRUE</code> then partial iteration details are displayed when <code>ASReml-R</code> functions are invoked; if <code>FALSE</code> then no output is displayed. |
| set.terms | a character vector specifying the terms that are to have constraints and/or initial values set prior to fitting. |
| ignore.suffices | a logical vector specifying whether the suffices of the <code>asreml</code> -assigned names of the variance terms (i.e. the information to the right of an "!", other than "R!") is to be ignored in matching elements of <code>terms</code> . If <code>TRUE</code> for an element of <code>terms</code> , the suffices are stripped from the <code>asreml</code> -assigned names. If <code>FALSE</code> for an element of <code>terms</code> , the element must exactly match an <code>asreml</code> -assigned name for a variance term. This vector must be of length one or the same length as <code>terms</code> . If it is of length one then the same action is applied to the <code>asreml</code> -assigned suffices for all the terms in <code>terms</code> . |
| constraints | a character vector specifying the constraints to be applied to the terms specified in <code>terms</code> . This vector must be of length one or the same length as <code>terms</code> . If it is of length one then the same constraint is applied to all the terms in <code>terms</code> . If any of the constraints are equal to <code>NA</code> then they are left unchanged for those terms. |
| initial.values | a character vector specifying the initial values for the terms specified in <code>terms</code> . This vector must be of length one or the same length as <code>terms</code> . If it is of length one then the same initial value is applied to all the terms in <code>terms</code> . If any of the <code>initial.values</code> are equal to <code>NA</code> then they are left unchanged for those terms. |
| ... | further arguments passed to <code>asreml</code> and to <code>wald.asreml</code> . |

Value

An `asrtests` object, which is a list containing:

1. `asreml.obj`: an `asreml` object containing the fit after the term has been omitted from the model;
2. `wald.tab`: a 4-column data.frame containing a pseudo-anova table for the fixed terms produced by `wald.asreml`;
3. `test.summary`: a data.frame with columns `term`, `DF`, `denDF`, `p` and `action`. A row is added to it for each term that is dropped, added or tested or a note that several terms have been

added or removed. A row contains the name of the term, the DF, the p-value and the action taken. Possible codes are: Dropped, Retained, Swapped, Unswapped, Significant, Nonsignificant, Absent, Added, Removed and Boundary. If the changed model did not converge, Unconverged will be added to the code. Note that the logical `asreml.obj$converge` also reflects whether there is convergence.

If the term is not in the model, then the supplied `asreml` object will be returned. Also, `reml.test` will have the likelihood ratio and the p-value set to NA and the degrees of freedom to zero. Similarly, the row of `test.summary` for the term will have its name, a p-value set to NA, and action set to Absent.

See Also

[asremlPlus-package](#), [asrtests](#), [newrcov.asrtests](#),
[choose.model.asrtests](#), [reml.lrt.asreml](#), [rmboundary.asrtests](#),
[newfit.asreml](#), [testswapran.asrtests](#), [addrm.terms.asrtests](#),
[sig.devn.reparam.asrtests](#)

Examples

```
## Not run:
data(Wheat.dat)
current.asr <- asreml(yield ~ Rep + WithinColPairs + Variety,
                     random = ~ Row + Column + units,
                     rcov = ~ ar1(Row):ar1(Column),
                     data=Wheat.dat)
current.asrt <- asrtests(current.asr, NULL, NULL)
current.asrt <- rmboundary.asrtests(current.asrt)
# Test Row autocorrelation
current.asrt <- testrcov.asrtests("~ Row:ar1(Column)", current.asrt,
                                label="Row autocorrelation", simpler=TRUE)

print(current.asrt)

## End(Not run)
```

| | |
|-----------------------------------|--|
| <code>testswapran.asrtests</code> | <i>Tests, using a REMLRT, whether the difference between current random model and one in which oldterms are dropped and newterms are added is significant.</i> |
|-----------------------------------|--|

Description

Fits a new random model using `asreml` by removing oldterms and adding newterms. If `simpler = FALSE` the model to be fitted must be more complex than the one whose fit has been stored in `asrtests.obj`. That is, the new model must have more parameters. However, if `simpler = TRUE` the model to be fitted must be simpler than the one whose fit has been stored in `asrtests.obj` in that it must have fewer parameters. The test is a REML ratio test that is performed using [reml.lrt.asreml](#), which is only valid if the models are nested. It compares the newly fitted model with the fit of the model in `asrttest.obj`. A row is added to the `test.summary` data.frame using the supplied label. If the newly fitted model is retained, any boundary terms are then removed using [rmboundary.asrtests](#).

Usage

```
testswapran.asrtests(oldterms = NULL, newterms = NULL, asrtests.obj,
  label = "Swap in random model", simpler = FALSE,
  alpha = 0.05, positive.zero = FALSE, denDF="default",
  trace = FALSE, update = TRUE,
  set.terms = NULL, ignore.suffices = TRUE,
  constraints = "P", initial.values = NA, ...)
```

Arguments

| | |
|-----------------|---|
| oldterms | terms, stored as a character, that are to be removed from the random model using asreml. |
| newterms | terms, stored as a character, that are to be added to the random model using asreml. |
| asrtests.obj | an asrtests object for a fitted model that is a list containing an asreml object, a wald.tab data.frame with 4 columns, and a data.frame with 5 columns that records any previous changes and tests in the fitted model. |
| simpler | a logical indicating whether the new model to be fitted. after the changes made as a result of swapping oldterms for newterms, is simpler than the already fitted model whose fit is stored in asrtests.obj. |
| alpha | the significance level for the test. |
| label | a character string to use as the label in test.summary and which indicates what is being tested. |
| positive.zero | a logical indicating whether the hypothesized values for the varaince components being tested are on the boundary of the parameter space. For example, this is be true for positively-constrained variance components that, under the reduced model, are zero. |
| denDF | Specifies the enthod to use in computing approximate denominator degrees of freedom when wald.asreml is called. Can be none to suppress the computations, numeric for numerical methods, algebraic for algebraic methods or default, the default, to autommatically choose numeric or algebraic computations depending on problem size. The denominator degrees of freedom are calculated according to Kenward and Roger (1997) for fixed terms in the dense part of the model. |
| trace | if TRUE then partial iteration details are displayed when ASReml-R functions are invoked; if FALSE then no output is displayed. |
| update | if TRUE then update.asreml is called to change the model. In doing this the arguments R.param and G.param are set to those in the asreml object stored in asrtests.obj so that the values from the previous model are used as starting values. If FALSE then a call is made to asreml in which the only changes from the previous call are that (i) models are modified for the supplied oldterms and newterms, and (ii) modifications specified via ... are made. |
| set.terms | a character vector specifying the terms that are to have constraints and/or initial values set prior to fitting. |
| ignore.suffices | a logical vector specifying whether the suffices of the asreml-assigned names of the variance terms (i.e. the information to the right of an "!", other than "R!") is to be ignored in matching elements of terms. If TRUE for an element of terms, the suffices are stripped from the asreml-assigned names. If FALSE |

| | |
|-----------------------------|---|
| | for an element of <code>terms</code> , the element must exactly match an <code>asreml</code> -assigned name for a variance term. This vector must be of length one or the same length as <code>terms</code> . If it is of length one then the same action is applied to the <code>asreml</code> -assigned suffices for all the terms in <code>terms</code> . |
| <code>constraints</code> | a character vector specifying the constraints to be applied to the terms specified in <code>terms</code> . This vector must be of length one or the same length as <code>terms</code> . If it is of length one then the same constraint is applied to all the terms in <code>terms</code> . If any of the constraints are equal to <code>NA</code> then they are left unchanged for those terms. |
| <code>initial.values</code> | a character vector specifying the initial values for the terms specified in <code>terms</code> . This vector must be of length one or the same length as <code>terms</code> . If it is of length one then the same initial value is applied to all the terms in <code>terms</code> . If any of the <code>initial.values</code> are equal to <code>NA</code> then they are left unchanged for those terms. |
| <code>...</code> | further arguments passed to <code>asreml</code> and <code>wald.asreml</code> . |

Value

An `asrtests` object, which is a list containing:

1. `asreml.obj`: an `asreml` object containing the fit after the term has been omitted from the model;
2. `wald.tab`: a 4-column data.frame containing a pseudo-anova table for the fixed terms produced by `wald.asreml`;
3. `test.summary`: a data.frame with columns `term`, `DF`, `denDF`, `p` and `action`. A row is added to it for each term that is dropped, added or tested or a note that several terms have been added or removed. A row contains the name of the term, the DF, the p-value and the action taken. Possible codes are: Dropped, Retained, Swapped, Unswapped, Significant, Nonsignificant, Absent, Added, Removed and Boundary. If the changed model did not converge, Unconverged will be added to the code. Note that the logical `asreml.obj$converge` also reflects whether there is convergence.

If the term is not in the model, then the supplied `asreml` object will be returned. Also, `reml.test` will have the likelihood ratio and the p-value set to `NA` and the degrees of freedom to zero. Similarly, the row of `test.summary` for the term will have its name, a p-value set to `NA`, and action set to `Absent`.

See Also

[asrtests](#), [choose.model.asrtests](#), [reml.lrt.asreml](#), [rmboundary.asrtests](#), [newfit.asreml](#), [testrcov.asrtests](#), [addrm.terms.asrtests](#), [sig.devn.reparam.asrtests](#)

Examples

```
## Not run:
current.asrt <- testswapran.asrtests(oldterms = "str(~ Cart/xDays, ~us(2):id(184))",
                                   newterms = "Cart/xDays",
                                   current.asrt, pos = FALSE,
                                   label = "Intercept/Slope correlation",
                                   simpler = TRUE)

print(current.asrt)

## End(Not run)
```

| | |
|-------------------|--|
| variofaces.asreml | <i>plot empirical variogram faces, including envelopes, as described by Stefanova, Smith & Cullis (2009)</i> |
|-------------------|--|

Description

A function that produces a plot for each face of an empirical 2D variogram based on residuals produced after the fitting of a model using the function `asreml`. It also adds envelopes to the plot by simulating data sets from a multivariate normal distribution with expectation equal to the fitted values obtained from the fixed and spline terms and variance matrix equal to the fitted variance matrix (Stefanova, Smith & Cullis, 2009). The plot is controlled by the `rcov` model, which must consist of two factors corresponding to the two physical dimensions underlying the data. It can also have a third term involving the `at` function that defines sections of the data, such as experiments in different environments. In this case, the two variogram faces are produced for each section.

Usage

```
variofaces.asreml(object, V, nsim=100, seed = NULL, tolerance=1E-10,
                  units = "ignore", update = TRUE, trace = FALSE,
                  graphics.device=NULL, ...)
```

Arguments

| | |
|------------------------------|---|
| <code>object</code> | An <code>asreml</code> object from a call to <code>asreml</code> in which the data argument has been set. |
| <code>V</code> | The fitted variance matrix, i.e. having the appropriate pattern and values given the model fitted to the observed data and the estimates of the parameters obtained. |
| <code>nsim</code> | The number of data sets to be simulated in obtaining the envelopes. |
| <code>seed</code> | A single value, interpreted as an integer, that specifies the starting value of the random number generator. |
| <code>tolerance</code> | The value such that eigenvalues less than it are considered to be zero. |
| <code>units</code> | A character indicating whether the BLUPs for <code>units</code> are added to the residuals when this reserved factor is included in the random model. Possible values are <code>addtoresiduals</code> and <code>ignore</code> . If standardized conditional residuals are plotted and the BLUPs for <code>units</code> are to be added then it is the standardized BLUPs that are added. |
| <code>update</code> | if <code>TRUE</code> then the arguments <code>R.param</code> and <code>G.param</code> are set to those in the <code>asreml</code> object supplied in <code>object</code> so that the values from the original model are used as starting values. If <code>FALSE</code> then calls are made to <code>asreml</code> in which the only changes from the previous call are (i) the model is fitted to simulated data and (ii) modifications specified via <code>...</code> are made, except that changes cannot be made to any of the models. |
| <code>trace</code> | if <code>TRUE</code> then partial iteration details are displayed when <code>ASReml-R</code> functions are invoked; if <code>FALSE</code> then no output is displayed. |
| <code>graphics.device</code> | A character specifying a graphics device for plotting. The default is <code>graphics.device = NULL</code> , which will result in plots being produced on the current graphics device. Setting it to <code>"windows"</code> , for example, will result in a <code>windows</code> graphics device being opened. |

... Other arguments that are passed down to the function `asreml`. Changes to the models are not allowed. Other changes are dangerous and generally should be avoided.

Details

The `rcov` model is scanned to ensure that it involves only two factors not included in the `at` function, and to see if it has a third factor in an `at` function. If so, the faces of the 2D variogram, each based on one of the two non-`at` factors, are derived from the residuals in the supplied `asreml` object using `asreml.variogram`, this yielding the observed variogram faces. If `aom` was set to `TRUE` for the `asreml` object, the standardized consitional residuals are used. Then `nsim` data sets are generated by adding the `fitted.values`, extracted from the `asreml` object, to a vector of values randomly generated from a normal distribution with expectation zero and variance matrix `V`. Each data set is analyzed using the model in object and the variogram values for the faces are obtained using `asreml.variogram` stored. Note, if the analysis for a data set does not converge in `maxiter` iterations, it is discarded and a replacement data set generated. The value of `maxiter` can be specified in the call to `variofaces.asreml`. Plots are produced for each face and include the observed values and the 2.5%, 50% & 97.5% quantiles.

Value

A list with the following components:

1. **face1**: a `data.frame` containing the variogram values on which the plot for the first dimension is based.
2. **face2**: a `data.frame` containing the variogram values on which the plot for the second dimension is based.

Author(s)

Chris Brien

References

Stefanova, K. T., Smith, A. B. & Cullis, B. R. (2009) Enhanced diagnostics for the spatial analysis of field trials. *Journal of Agricultural, Biological, and Environmental Statistics*, **14**, 392–410.

See Also

[asremlPlus-package](#), [asreml](#), [plotvariofaces.asreml](#), [simulate.asreml](#).

Examples

```
## Not run:
data(Wheat.dat)
current.asr <- asreml(yield ~ Rep + WithinColPairs + Variety,
                    random = ~ Row + Column + units,
                    rcov = ~ ar1(Row):ar1(Column),
                    data=Wheat.dat)
current.asrt <- asrtests(current.asr, NULL, NULL)
current.asrt <- rmboundary.asrtests(current.asrt)
# Form variance matrix based on estimated variance parameters
s2 <- current.asr$sigma2
gamma.Row <- current.asr$gammas[1]
gamma.unit <- current.asr$gammas[2]
```

```

rho.r <- current.asr$gammas[4]
rho.c <- current.asr$gammas[5]
row.ar1 <- mat.ar1(order=10, rho=rho.r)
col.ar1 <- mat.ar1(order=15, rho=rho.c)
V <- gamma.Row * fac.sumop(Wheat.dat$Row) +
  gamma.unit * diag(1, nrow=150, ncol=150) +
  mat.dirprod(col.ar1, row.ar1)
V <- s2*V

#Produce variogram faces plot (Stefanaova et al, 2009)
variofaces.asreml(current.asr, V=V)

## End(Not run)

```

Wheat.dat

Data for an experiment to investigate 25 varieties of wheat

Description

The data appears in Gilmour et al. [1995] and is from a field experiment designed to compare the performance of 25 varieties of wheat. An analysis of it using asreml is presented by Butler et al. (2010, Section 8.6), although they suggest that it is a wheat experiment.

The experiment was conducted at Slate Hall Farm, UK, in 1976 and was designed as a balanced lattice square with 6 replicates laid out in a 10×15 rectangular grid. The columns in the data frame are: Rep, Row, Column, WithinColPairs, Variety, yield. The response variable is the grain yield.

Usage

```
data(Wheat.dat)
```

Format

A data.frame containing 150 observations of 6 variables.

Source

Butler, D. G., et al. (2010). *Analysis of Mixed Models for S language environments: ASReml-R reference manual*. Brisbane, DPI Publications.

Gilmour, A. R., et al. (1995) Average Information REML: An efficient algorithm for variance parameter estimation in linear mixed models. *Biometrics*, **51**, 1440-1450.

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