Package `LMERConvenienceFunctions`

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Description The main function of the package is to perform backward selection of fixed effects, forward fitting of the random effects, and post-hoc analysis using parallel capabilities. Other functionality includes the computation of ANOVAs with upper- or lower-bound p-values and R-squared values for each model term, model criticism plots, data trimming on model residuals, and data visualization. The data to run examples is contained in package LCF_data.
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Description
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Details

\begin{verbatim}
Package: LMERConvenienceFunctions
Type: Package
Version: 2.10
Date: 2015-1-31
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LazyLoad: yes
\end{verbatim}

Author(s)
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Maintainer: Antoine Tremblay <trea26@gmail.com>

References


See Also


Examples

```r
## Not run:
#########################################################
# Load and format data.
#########################################################
library(LCFdata)
data(eeg)

# restrict to electrode Fz and 80–180 ms window
eeg <- eeg[eeg$Time >= 80 & eeg$Time <= 180,]
eeg <- eeg[,c("Subject", "Item", "Time", "Fz",
"Freq", "Length","WMC")]

# mean center FreqB
eeg$FreqBc <- eeg$FreqB - mean(eeg$FreqB)
# split FreqBc into 3 categories. Doesn't make sense,
# but it's merely for example

eeg$FreqBdc <- "high"
eeg$FreqBdc[eeg$FreqBc <= quantile(eeg$FreqBc)[3]] <- "mid"
eeg$FreqBdc[eeg$FreqBc <= quantile(eeg$FreqBc)[2]] <- "low"
eeg$FreqBdc <- as.factor(eeg$FreqBdc) 
eeg$FreqBdc <- relevel(eeg$FreqBdc, "low")

# mean center LengthB
```
eeg$L_{\text{LengthBc}} <- \text{eeg$L_{\text{LengthB}} - mean(eeg$L_{\text{LengthB}}}$

# mean center WMC
\text{eeg$WMCc} <- \text{eeg$WMC - mean(eeg$WMC)}$

# Demonstrate plotDensity3d.fnc. 
plotDensity3d.fnc(x = sort(unique(eeg$WMCc)),
y = sort(unique(eeg$L_{\text{LengthBc}})))

# Demonstrate plotRaw3d.fnc. 
plotRaw3d.fnc(data = eeg, response = "Fz", pred = "WMCc",
intr = "LengthBc", plot.type = "persp", theta = 150)

library(lme4)

# fit initial model
\text{m0} <- \text{lmer(Fz \sim \text{FreqBdc} + \text{LengthBc} + WMCc)^2 + (1 \mid \text{Subject}), data = eeg}
\text{m1} <- \text{lmer(Fz \sim \text{FreqBdc} + \text{LengthBc} + WMCc)^2 + (1 \mid \text{Subject}) +
(1 \mid \text{Item}), data = eeg)}

# which model to choose?
\text{relLik(m0, m1)}

# check model assumptions
\text{mcp.fnc(m1)}

# remove outliers
\text{eeg <- romr.fnc(m1, eeg, trim = 2.5)}
\text{eeg$n\text{.removed}}
\text{eeg$percent\text{.removed}}
\text{eeg <- eeg$data}

# update model
\text{m1} <- \text{lmer(Fz \sim \text{FreqBdc} + \text{LengthBc} + WMCc)^2 + (1 \mid \text{Subject}) +
(1 \mid \text{Item}), data = eeg)}

# re-check model assumptions
\text{mcp.fnc(m1)}

# forward-fit random effect structure (simple for the purposes
# of the example).
m2 <- ffranefLMER.fnc(model = m1, ran.effects =
c("(0 + LengthBc | Subject)", "(0 + WMCc | Item)"),
log.file = FALSE)

# backfit model m2. In this case, could use bffixefLMER_t.fnc instead.
m3 <- bffixefLMER_F.fnc(m2, log.file = FALSE)

# The calls to ffranefLMER.fnc and bffixefLMER_F.fnc could
# be replaced by a call to fitLMER.fnc. In this latter case, however,
# bffixefLMER_F.fnc would be called first, then the random effect
# structure would be forward fitted, and finally teh fixed effects
# would be backfitted again.
m3b <- fitLMER.fnc(model = m1, ran.effects = c("(0 + LengthBc | Subject)",
"(0 + WMCc | Item)"), backfit.on = "F", log.file = FALSE)
pamer.fnc(m3b)
# The results are the same. This may not necessarily be the case
# elsewhere. First forward fitting the random effect structure and
# then backfitting the fixed effects, potentially pruning irrelevant
# random effects, is probably the best approach. Nonetheless, there is
# no hard evidence to this effect.

# check model assumptions
mcp.fnc(m3)

# check significance of model terms
pamer.fnc(m3)

# Demonstrate mcposthoc.fnc and summary.mcposthoc.

# Only the intercept is significant. For purposes of the
# example, let's perform a posthoc analysis on FreqBdc on
# model m2.
m2.ph <- mcposthoc.fnc(model = m2, var = list(ph1 = "FreqBdc"))

# Now check if and how the different levels differ between
# each other. First check high vs mid and high vs low:
summary(m2.ph, term = "FreqBdchigh")
# Then low vs mid (the low vs high row is redundant from the
# above summary):
summary(m2.ph, term = "FreqBdclow")
# Note that none of the levels differ from each other. Indeed,
# the backfitting process indicated that the model only has an
# intercept (i.e., the FreqBc factor variable was not significant).

# Just to show how one would look at posthocs for interactions. Let's
# look at the effect of Length at each FreqB bin:
summary(object = m2.ph, term = "LengthBc")
# Does Length effect different Freq bins? Start with low
# versus mid and high
smry <- summary(object = m2.ph, term = "FreqBdclow:LengthBc")
# then mid versus low and high
smry <- summary(object = m2.ph, term = "FreqBdcmid:LengthBc")

# Demonstrate 'revived' version of #
# plotLMER.fnc and plotLMER3d.fnc. #
# Generate plot for Length X Freq with function plotLMER.fnc.
plotLMER.fnc(m2, pred = "LengthBc", intr = list("FreqBdc",
levels(eeg$FreqBdc), "beg", list(1 : 3, 1 : 3)))

# Plotting the Length:WMC interaction with plotLMER3d.fnc. It'll #
take a little bit of time.
plotLMER3d.fnc(m2,"LengthBc","WMCc")
# Plot it a second time to demonstrate caching. You can notice the #
speed up.
plotLMER3d.fnc(m2,"LengthBc","WMCc")

# Demonstrate modeling and #
# backfitting of glmer. #
# Split FreqBc into 2 categories.
eeg$FreqBdc <- "high"
eeg$FreqBdc[eeg$FreqBc<=median(eeg$FreqBc)] <- "low"
eeg$FreqBdc <- as.factor(eeg$FreqBdc)
eeg$FreqBdc <- relevel(eeg$FreqBdc, "low")

# Fit glmer model.
m4 <- glmer(FreqBdc ~ (Fz + LengthBc + WMCC)^2 + (1 | Subject),
family = "binomial", data = eeg)
summary(m4)

# Back fit fixed effects, forward fit random effects, and then #
# re-back fit fixed effects. Need to set argument backfit.on to "t".
m5 <- fitLMER.fnc(model = m4, ran.effects = "(0 + LengthBc | Subject)",
backfit.on = "t", log.file = FALSE)
summary(m5)

# Plot the 2-way interaction.
plotLMER.fnc(m5, pred = "Fz", intr = list("LengthBc",
quantile(eeg$LengthBc), "med",list(1:5,1:5)))

# Look at the same plot, but in 3d.
plotLMER3d.fnc(m5, pred = "Fz", intr = "LengthBc")

# Test backfitting on AIC, #
# BIC, llrt, relLik.AIC, and #
# relLik.BIC. #
# AIC
m.test <- bffixefLMER_F.fnc(m2, method = "AIC", log.file = FALSE)
m.test <- bffixefLMER_t.fnc(m2, method = "AIC", log.file = FALSE)
m.test <- bffixefLMER_t.fnc(m4, method = "AIC", log.file = FALSE)
m.test <- bffixefLMER_F.fnc(m4, method = "AIC", log.file = FALSE)

# BIC
m.test <- bffixefLMER_F.fnc(m2, method = "BIC", log.file = FALSE)
m.test <- bffixefLMER_t.fnc(m2, method = "BIC", log.file = FALSE)
m.test <- bffixefLMER_t.fnc(m4, method = "BIC", log.file = FALSE)
m.test <- bffixefLMER_F.fnc(m4, method = "BIC", log.file = FALSE)

# llrt
m.test <- bffixefLMER_F.fnc(m2, method = "llrt", log.file = FALSE)
m.test <- bffixefLMER_t.fnc(m2, method = "llrt", log.file = FALSE)
m.test <- bffixefLMER_t.fnc(m4, method = "llrt", log.file = FALSE)
m.test <- bffixefLMER_F.fnc(m4, method = "llrt", log.file = FALSE)

# rellik.AIC
m.test <- bffixefLMER_F.fnc(m2, method = "rellik.AIC", log.file = FALSE)
m.test <- bffixefLMER_t.fnc(m2, method = "rellik.AIC", log.file = FALSE)
m.test <- bffixefLMER_t.fnc(m4, method = "rellik.AIC", log.file = FALSE)
m.test <- bffixefLMER_F.fnc(m4, method = "rellik.AIC", log.file = FALSE)

# rellik.BIC
m.test <- bffixefLMER_F.fnc(m2, method = "rellik.BIC", log.file = FALSE)
m.test <- bffixefLMER_t.fnc(m2, method = "rellik.BIC", log.file = FALSE)
m.test <- bffixefLMER_t.fnc(m4, method = "rellik.BIC", log.file = FALSE)
m.test <- bffixefLMER_F.fnc(m4, method = "rellik.BIC", log.file = FALSE)

## End(Not run)

### Description

This function back-fits an initial LMER model either on upper- or lower-bound p-values obtained from function pamert.fnc, log-likelihood ratio testing (LLRT), AIC, BIC, rellik.AIC, or rellik.BIC.
Note that this function CANNOT be used with generalized linear mixed-effects models (glmers).

Usage

```r
bfFixefLMER_F.fnc(model, item = FALSE, method = c("F", "llrt", "AIC", "BIC", "relLik.AIC", "relLik.BIC"), threshold = NULL, alpha = NULL, alpha.item = NULL, prune.ranefs = TRUE, p.value = "upper", set.REML.FALSE = TRUE, keep.single.factors = FALSE, reset.REML.TRUE = TRUE, log.file = NULL)
```

Arguments

- `model`: A mer object (fitted by function lmer). Note that this function cannot be used with generalized linear mixed-effects models (glmers).
- `item`: Whether or not to evaluate the addition of by-item random intercepts to the model, evaluated by way of log-likelihood ratio test. Either FALSE (the default) or the column name (quoted) of the item identifier (e.g., "Item", or "Word").
- `method`: Backfitting method. One of "F" (p-value), "llrt", "AIC", "BIC", "relLik.AIC", or "relLik.BIC" (relative likelihood, see function rellik). Defaults to F. You can find information regarding differences between AIC and BIC from http://methodology.psu.edu/eres.
- `threshold`: Method-specific threshold for parameter selection. It refers to alpha in the case of "F" and "llrt", to the minimum reduction in likelihood in the case of "AIC" and "BIC", or to the minimum difference in probability in the case of "relLik.AIC" and "relLik.BIC". Defaults NULL, which means 0.05 for "F" and "llrt", 5 for "AIC" and "BIC", and 4 for "relLik.AIC" and "relLik.BIC".
- `alpha`: If the method is F, it is the p-value (from pamer.fnc) above which a model term is dropped. In this case, it defaults to the value passed to argument threshold, i.e., 0.05. Otherwise it is the p-value threshold above which a test (see method) is performed between a model with the term under consideration and a simpler model without it (in this case, defaults to 0, i.e. all terms will be tested).
- `alpha.item`: Alpha value for the evaluation of by-item random intercepts. Defaults to 0.05 or to the specified threshold.
- `prune.ranefs`: Logical. Whether to remove any random effect for which its variable is not also present in the fixed effects structure (with the exception of the grouping variables such as "Subjects" and "Items"). Defaults to TRUE. For example, if the random effects structure contains the terms Condition + ROI + Group, and the random effects structure contains the terms (1 | Subject) + (0 + TrialNum | Subject), the random effect (0 + TrialNum | Subject) will be pruned from the model given that it is not in the model's fixed effects structure.
- `p.value`: If method = "F", whether to use upper-bound ("upper"; the default) or lower-bound ("lower") p-values during backfitting.
- `set.REML.FALSE`: Logical. Whether or not to set REML to FALSE. Defaults to TRUE.
- `keep.single.factors`: Logical. Whether or not main effects are kept (not subjected to testing and reduction). Defaults to FALSE.
reset.REML.TRUE Logical. Whether or not to re-set the back-fitted model to REML = TRUE.

log.file Whether a back-fitting log should be saved. Defaults to NULL, which means that a log is saved in a temporary folder with the file name file.path(tempdir(), paste("bfixefLMER_F_log", date(), "txt"), sep = "_"). The path and file name of the log can be changed to whatever the use wishes. Set to FALSE to disable.

Details

The back-fitting process works as follows:

1. If argument method is not set to F, REML is set to FALSE;

2. First consider only highest-order interaction model terms:

   (a) If method is F, the model term with the highest ANOVA p-value is identified. If this p-value is higher than alpha, the model term is removed and a new model is fitted. This is repeated for each model term that has a p-value higher than the alpha value. The algorithm then moves on to step (b). If method is not F, the model term with the lowest p-value is identified and the following is evaluated:

      i. A new model without this model term is fitted;

      ii. The more complex and simpler models are compared by way of a log-likelihood ratio test in case method is "llrt", by way of AIC or BIC values in case method is "AIC" or "BIC", or by calculating the relLik based on AIC or BIC in case method is "relLik.AIC" or "relLik.BIC". If the result determines that the term under consideration does not increase model fit, it is removed; otherwise it is kept.

      iii. Move on to the next model term with the smallest p-value smaller than alpha and repeat steps (i)–(iii).

   (b) Once all highest-order interaction terms have been evaluated, go down to the second highest order interactions: Repeat steps (ai)–(aiii) with the following addition: If a term would be removed from the model, but it is part of a high-order interaction, keep it. Once all terms of the interaction level have been evaluated, move down to the next lower-order level until main effects have been evaluated, after which the process stops. If keep.single factors = TRUE, the process stops after the evaluation of all interaction terms.

3. If argument method is set to something else other than "F", set reset.REML.TRUE to TRUE (default) unless otherwise specified.

In brief, if method is set to "F", a term remains in the model if its p-value is equal to or greater than alpha; if method is set to something else, a term remains in the model if

   1. its p-value from the ANOVA is equal to or smaller than alpha;
   2. it significantly increases model fit as determined by the specified method;
   3. it is part of a significant higher-order interaction term.

This backfitting method was used in Newman, Tremblay, Nichols, Neville, and Ullman (2012). If factorial terms are included in the initial model, back-fitting on F is recommended.
Value

A mer model with back-fitted fixed effects is returned and a log of the back-fitting process is printed on screen and (by default) in a log file in a temporary file.

Warnings

Upper-bound $p$-values can be anti-conservative, while lower-bound $p$-values can be conservative. See function pamer.fnc.

Note

If you get this error:

```r
Error in model.frame.default(data = ..2, formula = log_Segment_Duration ~ :  
The ... list does not contain 2 elements
```

It is probably because you updated the model using function update and the data now appears as data = ..2 or something similar to this. You can check this by typing model@call. If this is the case, re-fit your model as lmer(DV ~ IV + IV + (ranef) data = dat).

Author(s)

Antoine Tremblay, Dalhousie University, <trea26@gmail.com> and Johannes Ransijn <johannesransijn@gmail.com>.

References


See Also

bfFixefLMER_t.fnc; ffRanefLMER.fnc; fitLMER.fnc; mcpPosthoc.fnc; pamer.fnc; mcp.fnc; relLik; romr.fnc

Examples

# see example in LMERConvenienceFunctions help page.
**Description**

This function back-fits an initial LMER model on $t$-values, and, if enabled, log-likelihood ratio testing. Note that, this function CAN be used with generalized linear mixed-effects models (glmers).

**Usage**

```r
bfFixefLMER_tNfnc(model, item = FALSE, method = c("t", "z", "llrt", "AIC", "BIC", "rellik.AIC", "rellik.BIC"), threshold = NULL, t.threshold = NULL, alpha.item = NULL, prune.ranefs = TRUE, set.REML.FALSE = TRUE, keep.single.factors = FALSE, reset.REML.TRUE = TRUE, log.file = NULL)
```

**Arguments**

- **model**: A mer object (fitted by function lmer). Note that this function can be used with generalized linear mixed-effects models (glmers).
- **item**: Whether or not to evaluate the addition of by-item random intercepts to the model, evaluated by way of log-likelihood ratio testing. Either FALSE (the default) or the column name (quoted) of the item identifier (e.g., "Item", or "Word").
- **method**: Backfitting method. One of "t" (lmer), "z" (glmer), "llrt", "AIC", "BIC", "rellik.AIC", or "rellik.BIC" (the latter two are based on relative likelihood, see function rellik). Defaults to "t". You can find information regarding differences between AIC and BIC from [http://methodology.psu.edu/eresources/ask/sp07](http://methodology.psu.edu/eresources/ask/sp07).
- **threshold**: Method-specific threshold for parameter selection. It refers to the minimum $t/z$-value in the case of "t" or "z", to the alpha value in the case of "llrt", to the minimum reduction in likelihood in the case of "AIC" and "BIC", or to the minimum difference in probability in the case of "rellik.AIC" and "rellik.BIC". Defaults NULL, which means 2 for "t" and "z", 0.05 for "llrt", 5 for "AIC" and "BIC", and 4 for "rellik.AIC" and "rellik.BIC".
- **t.threshold**: Defaults to NULL. If the method = "t" or method = "z", it is the $t/z$-value below which a model term is dropped (if t.threshold = NULL, it will be set to 2). Otherwise it is the threshold for $t/z$-value below which a test (see method) is performed between a model with the term under consideration and a simpler model without it (if t.threshold = NULL, it is set to Inf, which means that all terms are tested).
- **alpha.item**: Alpha value for the evaluation of by-item random intercepts. Defaults to 0.05 or to the specified threshold in case method is "llrt".
prune.ranefs Logical. Whether to remove any random effect for which its variable is not also present in the fixed effects structure (with the exception of the grouping variables such as "Subjects" and "Items"). Defaults to TRUE. For example, if the random effects structure contains the terms Condition + ROI + Group, and the random effects structure contains the terms (1 | Subject) + (0 + TrialNum | Subject), the random effect (0 + TrialNum | Subject) will be pruned from the model given that it is not in the model’s fixed effects structure.

set.REML.FALSE Logical. Whether or not to set REML to FALSE. Defaults to TRUE. Not used for glmer models.

reset.REML.TRUE Logical. Whether or not to re-set the back-fitted model to REML = TRUE. Not used for glmer models.

keep.single.factors Logical. Whether or not main effects are kept (not subjected to testing and reduction). Defaults to FALSE.

log.file Whether a back-fitting log should be saved. Defaults to NULL, which means that a log is saved in a temporary folder with the file name file.path(tempdir()), paste("bfFixefLMER_F", date(), "txt"), The path and file name of the log can be changed to whatever the use wishes. Set to FALSE to disable.

Details

The back-fitting process works as follows:

1. If argument method is not set to "t", REML is set to FALSE;
2. First consider only highest-order interaction model terms:
   (a) If method is "t" or "z", the model term with the lowest \( t/z \)-value is identified. If this \( t/z \)-value is smaller than threshold, the model term is removed and a new model is fitted. This is repeated for each model term for term that has a \( t \)-value smaller than the threshold value. The algorithm then moves on to step (b). If method is not "t" or "z", the model term with the lowest \( t/z \)-value-value-value is identified and the following is evaluated:
      i. A new model without this model term is fitted;
      ii. The more complex and simpler models are compared by way of a log-likelihood ratio test in case method is "llrt", by way of AIC or BIC comparison if method is "AIC" "BIC", or by calculating the relLik based on AIC or BIC in case method is "relLik.AIC" or "relLik.BIC". If the result determines that the term under consideration does not increase model fit, it is removed; otherwise it is kept.
      iii. Move on to the next model term with the smallest \( t/z \)-value smaller than threshold and repeat steps (i)–(iii).
   (b) Once all highest-order interaction terms have been evaluated, go down to the second highest order interactions: Repeat steps (a)–(a(iii)) with the following addition: If a term would be removed from the model, but it is part of a high-order interaction, keep it. Once all terms of the interaction level have been evaluated, move down to the next lower-order level until main effects have been evaluated, after which the process stops. If keep.single.factors = TRUE, the process stops after the evaluation of all interaction terms.
bfFixefLMER_t.fnc

3. If argument method is set to something other than t or z, set reset.REML.TRUE to TRUE (default) unless otherwise specified.

In brief, if method is set to "t" or "z", a term remains in the model if its $t/z$-value is equal to or greater than threshold; if method is set to something else, a term remains in the model if

1. its $t/z$-value is equal to or greater than threshold;
2. it significantly increases model fit as determined by the specified method;
3. it is part of a significant interaction term.

This backfitting method was used in Tremblay & Tucker (2011). If factorial terms with more than two levels are included in the initial model, back-fitting on F is recommended.

Value

A mer model with back-fitted fixed effects (on $t$-values) is returned and a log of the back-fitting process is printed on screen and (by default) in a log file.

Note

If you get this error:

```r
Error in model.frame.default(data = ..2, formula = log_Segment_Duration ~ : The ... list does not contain 2 elements
```

It is probably because you updated the model using function update and the data now appears as 
`data = ..2` or something similar to this. You can check this by typing `model@call`. If this is the case, re-fit your model as `lmer(DV ~ IV + IV + (ranef ~ data) ~ dat)`.

Author(s)

Antoine Tremblay, Dalhousie University, <trea26@gmail.com> and Johannes Ransijn <johannesransijn@gmail.com>.

References


See Also

`bfFixefLMER_F.fnc`; `ffRanefLMER.fnc`; `fitLMER.fnc`; `mcposthoc.fnc`; `pamer.fnc`; `mcp.fnc`; `rellik`; `romr.fnc`

Examples

```r
# see example in LMERConvenienceFunctions help page.
```
cd      

Change directory.

Description
Change directory to the one corresponding to the row number listed by function f.

Usage

cd(dir)

Arguments

dir The row number corresponding to the directory list returned by function f.

Value
Change directory to the selected one.

Author(s)
Antoine Tremblay, Dalhousie University, <trea26@gmail.com>

See Also
f; cdf; cdup; setwd

cdf     

Change directory; list files and directories in new directory using function f.

Description
Change directory to the one corresponding to the row number returned by function f.

Usage

cdf(dir)

Arguments

dir The row number corresponding to the directory listed by function f.

Value
Change to new directory and list files and directories in new directory using function f.
cdup

Author(s)
Antoine Tremblay, Dalhousie University, <trea26@gmail.com>

See Also
f; cd; cdup; setwd

cdup

Change directory one level up.

Description
Change directory one level up and list directory and files in new directory.

Usage
cdup()

Value
Change directory one level up.

Author(s)
Antoine Tremblay, Dalhousie University, <trea26@gmail.com>

See Also
f; cd; cdf; setwd

cn
List the column names of a data frame in matrix format.

Description
The column names of the specified data frame are listed in matrix format, that is, each one appears in one row preceded by the row number.

Usage
cn(data.frame)

Arguments
data.frame A data frame.
Value

A matrix containing the column names of the data frame.

Author(s)

Antoine Tremblay, Dalhousie University, <trea26@gmail.com>

See Also

colnames

f

List files and directories in current directory.

Description

List files and directories in current directory in matrix format. Each row is preceded by a row number.

Usage

f(path = ".", pattern = NULL, all.files = FALSE, full.names = FALSE, recursive = FALSE, ignore.case = FALSE)

Arguments

path A character vector of full path names; the default corresponds to the working directory getwd(). Missing values will be ignored.

pattern An optional regular expression. Only file names which match the regular expression will be returned.

all.files Logical. If FALSE, only the names of visible files are returned. If TRUE, all file names will be returned.

full.names Logical. If TRUE, the directory path is prepended to the file names. If FALSE, only the file names are returned.

recursive Logical. Should the listing recurse into directories?

ignore.case Logical. Should pattern-matching be case-insensitive?

Value

A matrix containing the names of the files and directories, preceded by a row number, in the specified directories. If a path does not exist or is not a directory or is unreadable it is skipped, with a warning.

The files are sorted in alphabetical order, on the full path if full.names = TRUE. Directories are included only if recursive = FALSE.
Note
File naming conventions are platform dependent. recursive = TRUE is not supported on all platforms and may be ignored (with a warning).

Author(s)
Antoine Tremblay, Dalhousie University, <trea26@gmail.com>

See Also
list.files

Examples
f()

ffRanefLMER.fnc  Forward-fit the random effect structure of an LMER model.

Description
Forward-fit an LMER model’s random effect structure by comparing a model without one of the specified random effects and a model with it by way of log-likelihood ratio testing. If the more complex model is a significantly better fit, the random effect is kept, otherwise it is dropped. This function can now be used with generalized linear mixed-effects models (glmers).

Usage
ffRanefLMER.fnc(model, ran.effects = list(ran.intercepts = as.character(), slopes = as.character(), corr = as.character(), by.vars = as.character()), alpha = 0.05, if.warn.not.add = TRUE, log.file = NULL)

Arguments
model  A mer object (fitted by function lmer). This function can now be used with generalized linear mixed-effects models (glmers).
ran.effects  Can be either a vector or a list. In the former case, the random effects to be evaluated are provided. For example c("(1 + Frequency | Subject)", "(0 + Length | Subject)", "(1 | Subject)"). In the latter case, the list can be composed of (i) a vector of random intercepts to be evaluated (ran.intercepts), (ii) a vector of random slopes to be evaluated (slopes), (iii) a vector specifying, for each element of slopes, whether the correlation between the slope and by-variables specified in by.vars should be added (corr), and (iv) a vector of "by" variables for the random slopes (by.vars). Values that can be supplied to the corr argument are 1 (add correlation), 0 (do not add correlation), and NA (for when the "slope" is a factor variable). Note that if a term in slopes is a factor variable, the corr value tied to
it will be automatically set to NA. Also note that if no values are supplied to \texttt{corr}, a vector of 0 as long as the \texttt{slopes} vector will be automatically supplied. For example \texttt{list(ran.intercepts = "Word", slopes = c("Frequency", "Length", "NSynSet", "Class")})

Another example is \texttt{list(slopes = c("Trial", "Class"), by.vars = "Subject")}, where the \texttt{corr} argument will be equal to \texttt{c(0, NA)}.

\begin{itemize}
    \item \textbf{alpha} Level of significance for log-likelihood ratio test. Defaults to 0.05.
    \item \textbf{if.warn.not.add} Logical. If a warning is issued after fitting a model with a new random effect (e.g., \texttt{false convergence} or the like), should the random effect nevertheless be evaluated? Defaults to \texttt{TRUE}, meaning that if such a warning is issued, the random effect will not be added to the random effects structure of the model. If set to \texttt{FALSE}, the random effect will be evaluated for inclusion as any other random effects would be via log likelihood ratio testing even if a warning is issued.
    \item \textbf{log.file} Should the back-fitting log be saved? Defaults to \texttt{NULL}, which means that a log file is saved in a temporary folder as \texttt{paste("ffRanefLMER_log", gsub("::", ",", gsub(" ", "_", c("", "", "", "", "", "", "", "", "", "", "", "", "", "", "", "", "", "", "", "", "", "", "", "", "", "", "", "", "", ""))))}. The path and file name of the log can be changed to whatever the user wishes. Set to \texttt{FALSE} to disable.
\end{itemize}

\textbf{Value}

A \texttt{mer} object with forward-fitted random effect structure as well as a log of the process is printed on screen and, optionally, printed in a log file.

\textbf{Note}

The removal of a random effect from the random effects structure if the variables that compose it are not also in the fixed effects structure has been turned off in this version.

\textbf{Author(s)}

Antoine Tremblay, Dalhousie University, trea26@gmail.com.

\textbf{References}


\textbf{See Also}

\texttt{bfinancialLMER.fnc; bffixefLMER_t.fnc; fitLMER.fnc; mcposthoc.fnc; pamer.fnc; mcp.fnc; rom.fnc; persubjecttrim.fnc;}

\textbf{Examples}

\texttt{
# see example in LMERConvenienceFunctions help page.}

Back-fit fixed effects and forward-fit random effects of an LMER model.

**Description**

The function follows these steps: (1) If llrt is set to TRUE, set REML to FALSE (unless specified otherwise); (2) back-fit initial model either on $F$- (by default) or on $t/z$-values; (3) forward-fit random effects; (4) re-back-fit fixed effects; (5) if llrt is set to TRUE, set REML to TRUE (unless specified otherwise). Note that, this function CAN be used with generalized linear mixed-effects models (glmers).

**Usage**

```r
fitlmerNfnc(modelL itemL FALSEL backfitNonL c(BfBL BtBL BzBL BllrtBL BaicBL BbicBL BrellikNaicBL BrellikNbicBL thresholdL nullL tNthresholdL nullL ranNeffectsL list(ran.interceptsL as.character(),
  slopesL as.character(), corrL as.character(),
  by.varsL as.character()), alphaL NULL, alphaitemL NULL,
  if.warn.not.addL TRUE, prune.ranefsL TRUE, p.valueL "upper",
  set.REML.FALSEL TRUE, keep.single.factorsL FALSE,
  reset.REML.TRUEL TRUE, log.file.nameL NULL)
```

**Arguments**

- **model**
  A `mer` object (fitted by function `lmer`). This function can be used with generalized linear mixed-effects models (glmers) if argument `backfit.on` is set to "t", but not if it is set to "F".

- **item**
  Whether or not to evaluate the addition of by-item random intercepts to the model, evaluated by way of log-likelihood ratio test. Either FALSE (the default, does not evaluate this addition) or the column name (quoted) of the item identifier (e.g., "Item", "Word").

- **backfit.on**
  Either "F" (default) or "t". Refers to the statistic which will be used to determine which term to test and potentially remove from the model. If you are backfitting a generalized linear mixed-effects model (glmer), make sure to set `backfit.on` to "t"; the algorithm effectively backfits on "$z$".

- **method**
  Backfitting method. One of "F" ($p$-value), "t" ($t$ statistic), "$z" ($z$ statistic), "llrt", "AIC", "BIC", "relLik.AIC", or "relLik.BIC" (the latter two are based on relative likelihood, see function `rellik`). Defaults to "t". You can find information regarding differences between AIC and BIC from [http://methodology.psu.edu/eresources/ask/sp07](http://methodology.psu.edu/eresources/ask/sp07).

- **threshold**
  Method-specific threshold for parameter selection. It refers to alpha in the case of "F" and "llrt", to the $t/z$-value in case of "t" or "$z" to the minimum reduction in likelihood in the case of "AIC" and "BIC", or to the minimum difference in probability in the case of "relLik.AIC" and "relLik.BIC". Defaults NULL, which means 0.05 for "F" and "llrt", 2 for "t", 5 for "AIC" and "BIC", and 4 for "relLik.AIC" and "relLik.BIC".
t.threshold

Defaults to NULL. If the method = "t" or method = "z", it is the \( t/z \)-value below which a model term is dropped (if \( t \).threshold = NULL, it will be set to 2). Otherwise it is the threshold for \( t/z \)-value below which a test (see method) is performed between a model with the term under consideration and a simpler model without it (if \( t \).threshold = NULL, it is set to Inf, which means that all terms are tested.

ran.effects

Can be either a vector or a list. In the former case, the random effects to be evaluated are provided. For example c("(1 + Frequency | Subject)", "(0 + Length | Subject)", "(1 + Score | Subject)"). In the latter case, the list can be composed of (i) a vector of random intercepts to be evaluated (ran.intercepts), (ii) a vector of random slopes to be evaluated (slopes), (iii) a vector specifying, for each element of slopes, whether the correlation between the slope and by-variables specified in by.vars should be added (corr), and (iv) a vector of "by" variables for the random slopes (by.vars). Values that can be supplied to the corr argument are 1 (add correlation), 0 (do not add correlation), and NA (for when the "slope" is a factor variable). Note that if a term in slopes is a factor variable, the corr value tied to it will be automatically set to NA. Also note that if no values are supplied to corr, a vector of 0 as long as the slopes vector will be automatically supplied. For example list(ran.intercepts = "Word", slopes = c("Frequency", "Length", "NSynSet","Class"), corr = c(1L, 0L, 1L), by.vars = "Subject"). Another example is list(slopes = c("Trial", "Class"), by.vars = "Subject"), where the corr argument will be equal to c(0L, NA).

alpha

If the method is F, it is the \( p \)-value (from pamer.fnc) above which a model term is dropped. In this case, it defaults to the value passed to argument threshold, i.e., 0.05. Otherwise it is the \( p \)-value threshold above which a test (see method) is performed between a model with the term under consideration and a simpler model without it (in this case, defaults to 0, i.e. all terms will be tested).

alpha.item

Alpha value for the evaluation of by-item random intercepts. Defaults to 0.05 or to the specified threshold.

if.warn.not.add

Logical. If a warning is issued after fitting a model with a new random effect (e.g., false convergence or the like), should the random effect nevertheless be evaluated? Defaults to TRUE, meaning that if such a warning is issued, the random effect will not be added to the random effects structure of the model. If set to FALSE, the random effect will be evaluated for inclusion as any other random effects would be via log likelihood ratio testing even if a warning is issued.

prune.ranefs

Logical. Whether to remove any random effect for which its variable is not also present in the fixed effects structure (with the exception of the grouping variables such as "Subjects" and "Items"). Defaults to TRUE. For example, if the random effects structure contains the terms Condition + ROI + Group, and the random effects structure contains the terms (1 | Subject) + (0 + TrialNum | Subject), the ranedom effect (0 + TrialNum | Subject) will be pruned from the model given that it is not in the model's fixed effects structure.

p.value

Whether to use upper-bound ("upper"; the default) or lower-bound ("lower") \( p \)-values when back-fitting with method "F".

set.REML.FALSE

Logical. Whether or not to set REML to FALSE. Defaults to FALSE.
reset.REML.TRUE

Logical. Whether or not to re-set the back-fitted model to REML = TRUE.

keep.single.factors

Logical. Whether or not main effects are kept (not subjected to testing and reduction). Defaults to FALSE.

log.file.name

Should the back-fitting log be saved? Defaults to NULL, which means that a log file is saved in a temporary folder (platform dependent) as file.path(tempdir(), paste("fitLMER_log", file.path(tempdir()), "temp_log", file.path(tempdir()), "")), The path and file name of the log can be changed to whatever the user wishes. Set to FALSE to disable.

Details

The process has three stages. In the first stage, either bffixefLMER_F.fnc or bffixefLMER_t.fnc is called (depending on the user’s choice) and the fixed effects are back-fitted accordingly. In the second stage, fffranefLMER.fnc is called and random effects are forward-fitted. In the third stage, the fixed effects are back-fitted again. This is done because the inclusion of certain random effects sometimes renders certain fixed effects non-significant. This process was used in Tremblay and Tucker (2011) and in Newman, Tremblay, Nichols, Neville, and Ullman (2012).

If, for example, you have many analyses to run and a cluster is available, write a bash script that will create (1) .R files that will relevel the conditions and update the model, and (2) an associated .sh job submission script to submit the .R files. For example, let’s consider two ERP analyses all in a time window ranging from 100 to 250 ms. Two three-way interactions were considered: Position (factor; 1 to 6) X Length of the second word of a four-word sequence (e.g., in the middle of) X Working Memory Capacity score (continuous, from 0 to 100) and Trial (continuous; 1 to 432) X Length X Working Memory Capacity. Analyses were performed at electrodes Fp1 Fp2 AF3 AF4 F7 F3 Fz F4 F8 FC5 FC1 FC2 FC6 T7 C3 C4 T8 CP5 CP1 CP2 CP6. See Tremblay and Newman (In preparation) for more details. The analysis script named Fp1–CP6_100250.sh we used on the ACEnet cluster is as follows:

electrodes=({Fp1 Fp2 AF3 AF4 F7 F3 Fz F4 F8 FC5 FC1 FC2 FC6 T7 C3 C4 T8 CP5 CP1 CP2 CP6})
for e in ${electrodes[*]}; do
  export E=$e;
  # create .R script to load data, perform necessary manipulations
  # and perform the analysis using fitLMER.fnc
  echo '<<-Sys.getenv("E")' >> $e".R"
  echo 'load("/data/eeg600_trim_v2.rda")' >> $e".R"
  echo 'dat0<dat' >> $e".R"
  echo 'rm(dat).gc(T,T)' >> $e".R"
  echo 'dat <- dat0[dat0$Time >= 100 & dat0$Time <= 250, , drop = TRUE]' >> $e".R"
  echo 'dat <- dat[dat$Electrode == e, , drop = TRUE]' >> $e".R"
  echo 'subj<sort(unique(dat$Subject))' >> $e".R"
  echo 'for(i in subj){' >> $e".R"
  echo 'tmp<dat[dat$Subject==i,drop=TRUE]' >> $e".R"
  echo 'tmp$newfact<paste(tmp$Block,tmp$Position,sep=" ")' >> $e".R"
  echo 'newvec<vector("numeric")' >> $e".R"
  echo 'for(j in 1:length(unique(tmp$newfact))){' >> $e".R"
  echo 'newvec<rep(j,nrow(tmp$newfact==unique(tmp$newfact)[j]))}' >> $e".R"
  echo '}' >> $e".R"
echo 'tmp$Trial<-newvec' >> $e".R"
echo 'if(grep(i,subj)[1]==1)' >> $e".R"
echo 'newdat<-tmp' >> $e".R"
else' >> $e".R"
else' <- rbind(newdat,tmp)' >> $e".R"
}
else' >> $e".R"
else' <- newdat' >> $e".R"
else' <- as.factor(dat$Position)' >> $e".R"
else' <- lmer(Amplitude ~ (Position + Trial)*(LengthBc + WMCc) + ') >> $e".R"
else' <- fitLMER.fnc(m7,item="Item",ran.effects= c("(0+Trial|Subject)"),' >> $e".R"
else' <- lmer(m7b,m7b,'(0+LengthBc|Subject)",(0+Trial|Item)",(0+WMCc|Item)")' >> $e".R"
else' <- smry<-pamer.fnc(m7b)'' >> $e".R"
else' <- save(m7b,file=file.path("..","models",paste("m7b_","e","_100250.rda",sep=""))') >> $e"
else' <- save(smry,file=file.path("..","summaries",paste("smry_m7b_","e","_100250.rda",sep=""))')

### create the job submission script for the .R file created above
### submit the job
psub "job."$e".sh"
done;

and then type in the console

. Fp1-CP6_100250.sh

On the ACEnet cluster, this results in 22 independent analyses, simultaneously using a total of 22 cores and 176 GB of RAM. This analysis completes in about 30 minutes to 1 hour.

Value

A mer object with back-fitted fixed effects and forward-fitted random effects, as well as a log of the process, which is printed on screen and, optionally, printed in a log file.

Warnings

Upper-bound p-values can be anti-conservative, while lower-bound p-values can be conservative. See function pamer.fnc.
Note

The removal of a random effect from the random effects structure if the variables that compose it are not also in the fixed effects structure has been turned off in this version.

Author(s)

Antoine Tremblay, Dalhousie University, <trea26@gmail.com>

References


See Also

bfixefLMER.F.fnc; bfixefLMER.t.fnc; ffranefLMER.fnc; mcposthoc.fnc; pamer.fnc; mcp.fnc; relLik;

Examples

# see example LMERConvenienceFunctions help page.

mcp.fnc

Model criticism plots.

Description

A function to graph criticism plots for an LMER model (as in Baayen, 2008, chapter 7). Note that this function cannot be used with generalized linear mixed-effects models (GLMERs). Also note that the fourth plot (dfits) is omitted until we can figure out how to calculate dfits for a merMod object.

Usage

mcp.fnc(model, trim = 2.5, col = "red")
Arguments

model  A mer object (fitted by function lmer). Note that, at the moment, this function cannot be used with generalized linear mixed-effects models (GLMERs).

trim  Used to plot lines in the fitted ~ standardized residuals plot. The lines correspond to the threshold at which residuals would be or were removed. Defaults to 2.5 (standard deviations above and below the residuals mean).

col  Color of the lines added to the quantile-quantile plot and fitted ~ standardized residuals plot. Defaults to red.

Details

The first of the four plots graphs the density of the model residuals. The second plot graphs the quantile-quantile plot (actual standardized residuals versus theoretical quantiles). The third plot illustrates the fitted values versus the standardized residuals. The fourth graph plots the absolute values of the dffits of the residuals (not producing this plot as of version 2.2; might come back in future versions).

Value

Returns the four plots described above.

Author(s)

Antoine Tremblay, Dalhousie University, <trea26@gmail.com>.

References


Examples

# see example LMERConvenienceFunctions help page.

---

**mcposthoc.fnc**  Posthoc analyses for LMER models using parallel capabilities.

Description

This function uses the parallel package. For each factor level, a slave process is sent to one of the computer’s cores using function mclapply where the specified factor variables are re-leveled to each one of their levels, the mer model updated, and summaries returned. **MCMC p-value calculation is now implemented.** R will wait until all slave processes have finished running. See package parallel for more information about parallel computing. Note that traditional sequential computing can be achieved by specifying mc.cores = 1. Posthoc results can be viewed with function summary.mcposthoc.
Usage

mcposthoc.fnc(model, var, two.tailed = TRUE,
              mcmc = FALSE, nsim = 10000, ndigits = 4, mc.cores = 1,
              verbosity = 1, ...)  

Arguments

model A `mer` object (fitted by function `lmer`) or an `lm` object (fitted by function `lm`).
var A named list of variable on which to perform the posthoc analysis. For example
      list(ph1 = c("PronomOfTheme", "AnimacyOfRec", "DefinOfRec"), ph2 = c("SemanticClass"))
two.tailed Logical. Whether to perform one- or two-tailed t-tests. Defaults to `TRUE`, i.e.,
two-tailed.
mcmc Logical. Whether to calculate p-values using function `pamer.fnc` (the default)
or using function `pvals.fnc` from package `languager`.
nsim An integer denoting the required number of Markov chain Monte Carlo samples.
      Defaults to 10000.
ndigits Integer indicating the number of decimal places to be used in the t tables. De-
     faults to 4.
mc.cores The number of cores to use, i.e. how many processes will be spawned (at most).
verbosity Numeric. The amount of information printed to screen during the modeling
      process. The higher the number, the more information is printed. 0 turns this
      option off. Defaults to 1.
... Further arguments to pass to "mclapply".

Details

If var = list(ph1 = c("PronomOfTheme", "AnimacyOfRec", "DefinOfRec")), for example, the function will re-level and update the model on each combination of the variable levels as follows:

1. data$PronomOfTheme <- relevel(data$PronomOfTheme = "nonpronominal")
   data$AnimacyOfTheme <- relevel(data$AnimacyOfTheme = "animate")
   data$DefinOfTheme <- relevel(data$DefinOfTheme = "definite")

2. data$PronomOfTheme <- relevel(data$PronomOfTheme = "nonpronominal")
   data$AnimacyOfTheme <- relevel(data$AnimacyOfTheme = "inanimate")
   data$DefinOfTheme <- relevel(data$DefinOfTheme = "definite")

3. data$PronomOfTheme <- relevel(data$PronomOfTheme = "nonpronominal")
   data$AnimacyOfTheme <- relevel(data$AnimacyOfTheme = "animate")
   data$DefinOfTheme <- relevel(data$DefinOfTheme = "indefinite")

4. data$PronomOfTheme <- relevel(data$PronomOfTheme = "pronominal")
   data$AnimacyOfTheme <- relevel(data$AnimacyOfTheme = "animate")
   data$DefinOfTheme <- relevel(data$DefinOfTheme = "definite")

5. data$PronomOfTheme <- relevel(data$PronomOfTheme = "nonpronominal")
On a cluster, instead of using mcposthoc.fnc it is better (faster and less complicated) to write a bash script that will create (1) .R files that will relevel the conditions and update the model, and (2) an associated .sh job submission script to submit the .R files. For example, let’s consider two ERP analyses (regular past tense inflection and phrase structure) with three time windows each (300–400 ms, 550–700 ms, 750–850 ms in the regular past tense analysis, and 300–400 ms, 400–600 ms, and 750–850 ms in the phrase structure analysis). We investigated the effects of proficiency on ERP amplitudes. The initial models included a four-way interaction between Region of Interest (ROI) – with levels left anterior, left central, left posterior, midline anterior, midline central, midline posterior, right anterior, right central, and right posterior – Group (with levels L1 and L2), Condition (with levels control and violation), and Proficiency. After back-fitting the fixed effects, forward-fitting random effects, and reback-fitting the fixed effects as per fitlmer.fnc, the four-way interaction remained in every model. See Newman et al. (in preparation) for more details.

The posthoc analysis script named posthocs.sh we used on the ACEnet cluster is as follows:

```bash
for t in $(time[*]); do for i in $(condition[*]); do for j in $(group[*]); do for k in $(roi[*]); do
    ### create .R file where the model is updated on the data where
    ### re-level on each possible combination of variable levels
    export CONDITION=$i;
    export GROUP=$j;
    export ROI=$k;
    echo 'condition<-Sys.getenv("CONDITION")' >> "ph"$t$CONDITION$GROUP$ROI".R"
    echo 'group<-Sys.getenv("GROUP")' >> "ph"$t$CONDITION$GROUP$ROI".R"
    echo 'roi<-Sys.getenv("ROI")' >> "ph"$t$CONDITION$GROUP$ROI".R"
    echo 'load("models/m1"$t'.rda')' >> "ph"$t$CONDITION$GROUP$ROI".R"
    echo 'dat<m1@frame' >> "ph"$t$CONDITION$GROUP$ROI".R"
    echo 'dat$Condition<-relevel(dat$Condition,\'condition\')' >> "ph"$t$CONDITION$GROUP$ROI".R"
    echo 'dat$Group<-relevel(dat$Group,\'group\')' >> "ph"$t$CONDITION$GROUP$ROI".R"
    echo 'dat$ROI<-relevel(dat$ROI,\'roi\')' >> "ph"$t$CONDITION$GROUP$ROI".R"
    echo 'm1<-update(m1,\'-\',data=dat)' >> "ph"$t$CONDITION$GROUP$ROI".R"
```
Echo 'save(m1,file="ph$CONDITION$GROUP$ROI'.rda")' >> "ph$CONDITION$GROUP$ROI".R

### create the job submission script for the .R file created above

Echo '#$ -S /bin/bash' > "job.ph$CONDITION$GROUP$ROI".sh
Echo '#$ -cwd' >> "job.ph$CONDITION$GROUP$ROI".sh
Echo '#$ -j y' >> "job.ph$CONDITION$GROUP$ROI".sh
Echo '#$ -l h_rt=48:00:00' >> "job.ph$CONDITION$GROUP$ROI".sh
Echo '#$ -l h_vmem=8G' >> "job.ph$CONDITION$GROUP$ROI".sh
Echo '#$ -R y' >> "job.ph$CONDITION$GROUP$ROI".sh
Echo '#$ -N "ph$CONDITION$GROUP$ROI"' >> "job.ph$CONDITION$GROUP$ROI".sh
Echo 'R -q -f ph$CONDITION$GROUP$ROI'.R' >> "job.ph$CONDITION$GROUP$ROI".sh

### submit the job

Qsub "job.ph$CONDITION$GROUP$ROI".sh
done; done; done; done

and then type in the console

.posthoc.sh

On the AEnet cluster, this results in 2 * 3 * 9 * 2 * 2 = 216 independent analyses, simultaneously using a total of 216 cores and 1728 GB of RAM. This posthoc analysis completes in about 3-6 hours.

**Value**

An object of class "mcposthoc" with the following slots:

- **n**: The number of data points in data frame data.
- **var**: A named list containing the names of the variables used in the posthoc.
- **summaries**: A named list containing the posthoc summaries for each factor re-leveling. If mcmc = FALSE, data frames with upper- and lower-bound (anti-conservative and conservative, respectively) dfs, p-values, and deviance explained (%) for each model term. If mcmc = TRUE, data frames with the estimated coefficients, their MCMC mean, the HPD 95 and the probability based on the t distribution with the number of observations minus the number of fixed-effects coefficients as degrees of freedom. This last p-value is anti-conservative, especially for small data sets.

**warning**

Parallel computing capabilities will not be available on Windows because mclapply relies on forking. Sequential computing, however, will work on Windows if mc.cores = 1 (the default).
Note

It is not possible anymore to get p-values with function pvals.fnc of package languageR. Please see http://stackoverflow.com/questions/19199713/lme4-and-language-r-compatibility-error-input-model-is for other possible avenues to get p-values.

Author(s)

Antoine Tremblay, Dalhousie University, <trea26@gmail.com>.

See Also

summary.mcposthoc

Examples

# see example in LMERConvenienceFunctions help page.

Description

Compute upper- and lower-bound p-values for the analysis of variance (or deviance) as well as the amount of deviance explained (%) for each fixed-effect of an LMER model. Note that, at the moment, this function cannot be used with generalized linear mixed-effects models (glmers).

Usage

pamer.fnc(model, ndigits = 4)

Arguments

model A mer object (fitted by function lmer). Note that, at the moment, this function cannot be used with generalized linear mixed-effects models (glmers).

ndigits Integer indicating the number of decimal places to be used in the ANOVA table.

Details

Upper-bound p-values are computed by using as denominator \( df = nrow(model@frame) - qr(model@X)4rank \) (i.e., number of data points minus number of fixed effects including the intercept), which are anti-conservative. Lower-bound p-values are computed by using as denominator \( df = nrow(model@frame) - qr(model@X)4rank - \) (e.g., if by-subject intercepts and slopes, and there are 10 subjects, \( 10 * 2 = 20 \)). The amount of deviance explained by each model term (i.e., eta squared) is calculated as \([\text{Sum of Squares for the effect}] / [\text{Sum of Squares Total}]\). More specifically: \( \frac{\text{as.data.frame(anova(model))[2,2]}}{\text{sum((model@frame[, dv]-mean(model@frame[, dv]))^2)}} \) where dv is a vector of the names of the independent variables in the model.
Value

This function returns an object of class `data frame` with upper- and lower-bound (anti-conservative and conservative, respectively) `dfs`, `p-values`, and deviance explained (%) for each model term.

Author(s)

Antoine Tremblay, Dalhousie University, <trea26@gmail.com>

Examples

```r
# see example LMERConvenienceFunctions help page.
```

---

**perSubjectTrim.fnc**  
Per-subject Trimming of Response Variable.

---

Description

For each subject, removes data points that are, e.g., 2.5 standard deviations above or below the subject mean.

Usage

```r
perSubjectTrim.fnc(data, response, subject, trim = 2.5)
```

Arguments

- `data`: The data frame containing the data to be trimmed.
- `response`: The quoted name of the column containing the to-be-trimmed data.
- `subject`: The quoted name of the column containing subject identifiers.
- `trim`: Threshold at which data points will be removed. Defaults to 2.5 (standard deviations above and below each subject’s mean).

Value

The function returns the following objects:

- `data`: The data with outliers removed.
- `data0`: The original data prior to removing the outliers.
- `n.removed`: The number of data points removed.
- `percent.removed`: The percentage of removed data points.

Author(s)

Antoine Tremblay, Dalhousie University, <trea26@gmail.com>.
plotDensity3d.fnc

Kernel density estimation for two continuous variables.

Description

The densities of two continuous variables is first computed using the density function from package stats. The outer product of the two densities is computed, which can be plotted as a contour map, a perspective plot, or a dynamic 3d perspective graph.

Usage

plotDensity3d.fnc(x, y, plot.type = "contour", color = "terrain",
                 xlab = NULL, ylab = NULL, zlab = NULL, main = NULL, cex = 1,
                 alpha = 1, lit = TRUE, theta = 0, phi = 0, bw = "nrdf4",
                 adjust = 1, kernel = c("gaussian", "epanechnikov",
                 "rectangular", "triangular", "biweight", "cosine",
                 "optcosine"), weights = NULL, window = kernel, width, give.Rkern = FALSE,
                 n = 50, from, to, cut = 3, na.rm = FALSE, ...)

Arguments

x, y                  Numeric vectors.
plot.type             The type of plot to make. Can be any of "contour" (default), "persp", or, if package rgl is available, "persp3d".
color                 The colour scheme to use for plots. One of “topo”, “heat”, “cm”, “terrain”, “gray” or “bw”. Schemes “gray” and “bw” also modify the colors used.
xlab, ylab, zlab
   Titles for the axes. N.B. These must be character strings; expressions are not accepted. Numbers will be coerced to character strings.

main
   The main title on top of the plot.

cex
   The size of label and main text.

alpha
   For plot.type = "persp3d", alpha values between 0.0 (fully transparent) to 1.0 (opaque) for the main 3d surface.

lit
   Logical, specifying if lighting calculation should take place on geometry.

theta
   Angle defining the viewing direction. theta gives the azimuthal direction. Used only if plot.type = "persp".

phi
   Angle defining the viewing direction. phi gives the colatitude. Used only if plot.type = "persp".

bw, adjust, kernel, weights, window, width, give.Rkern, n, from, to, cut, na.rm
   See help page to function density.

... Further arguments passed to functions image, contour, persp, or persp3d.

Details

See help page to the density function as well as to Duncan Murdoch's persp3d function for more information. To save screenshots of "persp3d" plots (after plotting), use function rgl.snapshot (produces png files) or function rgl.postscript (produces eps files).

Value

Either a contour map or a (dynamic) perspective plot. Invisibly returns

x
   The numeric vector supplied in argument x.

y
   The numeric vector supplied in argument y.

xd
   The density object tied to vector x.

yd
   The density object tied to vector y.

mat
   The outer product of the x and y densities in matrix format.

col
   The color used for plotting.

Author(s)

Antoine Tremblay, Dalhousie University, <treaRV@gmail.com>.

See Also

contour; persp; density; outer

Examples

# see example in LMERConvenienceFunctions help page.
**Description**

Plot partial effects of a (generalized) linear mixed-effects model fit with lmer (compatible with package lme4 version > 1.0).

**Usage**

```r
plotLMER.fnc(model, xlabel = NA, xlabs = NA, ylabel = NA,
              ylimit = NA, ilabel = NA, fun = NA, pred = NA, control = NA,
              ranefs = NA, n = 100, intr = NA, lockYlim = TRUE, addlines = FALSE,
              withList = FALSE, cexsize = 0.5, linecolor = 1,
              addToExistingPlot = FALSE, verbose = TRUE, ...)
```

**Arguments**

- `model`: a mer model object
- `xlabel`: label for X-axis (if other than the variable name in the original model formula)
- `xlabs`: character vector with labels for X-axes in multipanel plot (if other than the variable names in the original model formula); if used, `xlabel` should not be specified
- `ylabel`: label for Y-axis (if other than the variable name of the dependent variable in the original model formula)
- `ylimit`: range for vertical axis; if not specified, this range will be chosen such that all data points across all subplots, including HPD intervals, will be accommodated
- `ilabel`: label for the interaction shown in the lower right-hand margin of the plot, overriding the original variable name in the model formula
- `fun`: a function to be applied for transforming the dependent variable, if NA, no transformation is applied; for models with `family = "binomial"`, `fun` is set to `plogis` by default; this can be disabled by setting `fun=function(x) return(x)`.
- `pred`: character string with name of predictor; if specified, a single plot will produced for the partial effect of this specific predictor
- `control`: a two-element list `list(predictor, val)` specifying a predictor the value of which has to be set to `val` in the partial effect plot(s); the predictor name should be exactly as specified in `names(model@fixef)`. It is up to the user to make sure that name and value make sense, the code here hands full 'control' to the user.
- `ranefs`: a four-element list `Group, Level, Predictor, specifying a random-effect Group (e.g. Subject), a level (e.g., S10) and a value (e.g., LogFrequency) for which partial effects have to be calibrated.
- `n`: integer denoting number of points for the plot, chosen at equally spaced intervals across the empirical range of the predictor variable
intr  

A list specifying an interaction to be graphed; obligatory arguments are (1) the name of the interaction variable, followed by (2) a vector of values for that variable, followed by (3) the position for interaction labels ("beg", "mid", or "end", or 'NA' if no labels are desired), optionally followed by (4) a list with as first element a vector of colors and as second element a vector of line types. The number of elements in both vectors should match the number of values specified under (2) for the interaction predictor.

lockylim  

Logical specifying whether all subplots should have the same range of values for the vertical axis; if TRUE, this range will be chosen to accommodate all fitted values including HDP intervals for all predictors across all plots.

addlines  

If TRUE, adds line(s) between levels of same factor(s).

withList  

Logical, if TRUE, a list will be output with all data frames for the subplots.

cexsize  

Character expansion size (cex) for additional information in the plot for interactions.

linecolor  

Color of lines in the plot, by default set to 1 (black).

addToExistingPlot  

Default FALSE, if set to TRUE, plot will be added to previous plot, but only if pred is specified.

verbose  

If TRUE (default), effect sizes and default transformations are reported.

...  

Further graphical parameters to be passed down; warning: col, pch, lty and cex will often generate an error as they are internally already fully specified for specialized subplots.

Details

When no predictor is specified, a series of plots is produced for the partial effects of each predictor. The graphs are shown for the reference level for factors and are adjusted for the median value for the other numerical predictors in the model. Interactions are not shown. The user should set up the appropriate number of subplots on the graphics device before running plotLMER.fnc().

Instead of showing all predictors jointly, plotLMER.fnc() can also be used to plot the partial effect of a specific predictor. When a specific predictor is specified (with pred = ...), a single plot is produced for that predictor. In this case, the intr argument can be used to specify a single second predictor that enters into an interaction with the selected main predictor.

Polynomials have to be fitted with poly(..., degree, raw=TRUE) and restricted cubic splines with rcs() from the rms package.

Note that any MCMC capabilities available in the languageR version of this function are not available in this version.

Value

A plot is produced on the graphical device.

Note

This code needs much more work, including (i) extension to poly with raw=FALSE, and (ii) general clean-up of the code.
Author(s)
R. H. Baayen, tweaked by Antoine Tremblay

See Also
plotLMER3d.fnc.

Examples
# see example in LMERConvenienceFunctions help page.

---

plotLMER3d.fnc  Dynamic 3d plot for mer object.

Description
Plot dynamic 3d partial effects of a (generalized) linear mixed-effects model fit with LMER.

Usage
plotLMER3d.fnc(model = NULL, pred, intr, plot.type = "contour",
xlim = range(x, na.rm = TRUE), ylim = range(y, na.rm = TRUE),
zlim = range(z, na.rm = TRUE), xlab = NULL,
ylab = NULL, zlab = NULL, main = NULL, shift = 0, scale = 1, cex = 1,
fun = NA, n = 30, color = "topo", alpha = 1, alpha.rs = 0.65, alpha.u = 1,
lit = TRUE, theta = 0, phi = 0, contourstepsize = 0.2, legend.args = NULL,
play3d = FALSE, ref.surf = FALSE, underneath = FALSE, add.raw = FALSE,
color.raw = "grey", alpha.raw = 0.5, rug = FALSE, rug.u = FALSE,
plot.dat = "default", path = "default", ...)

Arguments
model  A mer object or NULL (the default) to plot from an existing data-plotting object returned by this function and saved as an .rda file.
pred  The quoted name of a model predictor.
intr  The quoted name of a continuous model predictor.
plot.type  The type of plot to make. Can be any of "contour" (default), "image.plot" if package fields is available, "persp", or, if package rgl is available, "persp3d".
xlim, ylim, zlim  x-, y- and z-limits. The plot is produced so that the rectangular volume defined by these limits is visible.
xlab, ylab, zlab  Titles for the axes. N.B. These must be character strings; expressions are not accepted. Numbers will be coerced to character strings.
main  The main title on top of the plot.
shift
Constant to add to the smooth (on the scale of the linear predictor) before plotting. Defaults to 0. Passed to plotRaw3d.fnc.

scale
Constant by which to multiply the smooth before plotting. Defaults to 1. Passed to plotRaw3d.fnc.

cex
The size of label and main text.

fun
A function to be applied for transforming the dependent variable, if NA, no transformation is applied; for models with family = "binomial", fun is set to plogis by default; this can be disabled by setting fun=function(x)return(x).

n
Integer denoting number of points for the plot, chosen at equally spaced intervals across the empirical range of the predictor variable.

color
The colour scheme to use for plots. One of topo, heat, cm, terrain, gray or bw. Schemes gray, grey, and bw also modify the colors used.

alpha, alpha.rs, alpha.raw, alpha.u
For plot.type = "persp3d", alpha values between 0.0 (fully transparent) to 1.0 (opaque) for the main 3d surface, the reference surface, the added raw surface, and the "underneath" surface, respectively.

lit
Logical, specifying if lighting calculation should take place on geometry.

theta
Angle defining the viewing direction. theta gives the azimuthal direction. Used only if plot.type = "persp".

phi
Angle defining the viewing direction. phi gives the colatitude. Used only if plot.type = "persp".

contourstepsize
The size of the steps from contour line to contour line.

legend.args
When plot.type = "image.plot", arguments for a complete specification of the legend label. This is in the form of list and is just passed to the mtext function. (See example in image.plot help page). Defaults to NULL.

play3d
If plot.type = "persp3d" and play3d is set to TRUE, the 3d plot will spin around axis c(0, 0, 1) at rpm 4 for duration 20 seconds. The axis, rpm, and duration can be changed by supplying a three-argument list where the first argument is a three-element vector for the rotation axis, the second argument is an integer for the rotations per minute (rpm), and the the third argument is a rotation duration time.

ref.surf
If plot.type = "persp3d", whether a reference surface at the mean ought to be plotted. Defaults to FALSE.

underneath
If plot.type = "persp3d", whether a flat mirror image of the 3d surface ought to be plotted underneath it. Defaults to FALSE.

add.raw
If plot.type = "persp3d", whether to add a surface representing the raw data. Defaults to FALSE.

color.raw
The colour scheme to use for the raw data surface. One of topo, heat, cm, terrain, gray or bw. Schemes gray, grey, and bw also modify the colors used.

rug
Whether a rug ought to be plotted on the 3d surface. Defaults to FALSE.

rug.u
For plot.type = "persp3d", whether a rug ought to be plotted on the flat mirror image of the 3d surface. Defaults to FALSE.
Whether to cache the plotting data generated by a previous call to plotLMER3d.fnc. Generating the 3d plots can be time consuming. If the plot.dat argument is non-FALSE, the plotting information generated in the first call to the function will be saved so that in a second call to the function with exactly the same argument values, the plotting information will be retrieved and plotting will be significantly quicker. If plotting.data = "default" and path = "default", the plotting information will be saved in a temporary directory and the name of the file containing the information will equal to paste("lmer___", model@call, pred, intr, ".rda", sep = "."). The name of the file and the path where it will be saved can be set by the user in the plot.dat and path arguments. For example, plot.dat = "my_plotting.data", path = "Documents".

Note that "lmer___" will be appended to the beginning of whatever is specified in plot.dat and ".rda" to the end. Also note that if the user wants to save the plotting information returned by this function, the name of this object has to be z.

Further arguments to be passed to image, contour, image.plot, persp, or persp3d.

Details

See help page to Harald Baayen’s plotLMER.fnc function as well as to Duncan Murdoch’s persp3d function and the help page to function image.plot from package fields. To save screenshots of "persp3d" plots (after plotting), use function rgl.snapshot (produces png files) or function rgl.postscript (produces eps files).

Value

Invisibly returns plotting information (x and y vectors, z matrix, and colors, col). If plot.type = "contour", plot.type = "image.plot", or plot.type = "persp", a contour or perspective plot, respectively. If plot.type = "persp3d", a 2d plot as created by plotLMER.fnc as well as a dynamic 3d plot as created by persp3d. If ret = TRUE, a two-element list is returned containing the matrix and the matrix of corresponding colors is returned. If argument intel in non-null, a file containing plotting information will be saved.

Author(s)

Antoine Tremblay, Dalhousie University, <trea26@gmail.com>

See Also

persp; contour; plotLMER.fnc.

Examples

```r
if(try(require(LCFdata, quietly=TRUE))){
  data(z)
  temp.dir <- tempdir()
  save(z, file=file.path(temp.dir,"lmer___z.rda"))

  plotLMER3d.fnc(pred = "LengthBc", intr = "WMCC", plot.dat = "z", path = temp.dir)
```
plotLMER3d.fnc(pred = "Lengthbc", intr = "WMCc",
    plot.type = "persp", phi = 25, plot.dat = "z",
    path = temp.dir)
if(try(require(rgl, quietly=TRUE))){
    require(rgl)
    open3d()
    plotLMER3d.fnc(pred = "Lengthbc", intr = "WMCc",
        plot.type = "persp3d", plot.dat = "z", path = temp.dir)
}

plotRaw3d.fnc Visualize raw surface averages (3d)

Description
For a specified response variable and interacting continuous predictors, visualize in 3d the surface average.

Usage
plotRaw3d.fnc(data = NULL, response = NULL, pred = NULL, intr = NULL,
    xy = TRUE, color = "topo", zlim = NULL, xlab = NULL, ylab = NULL,
    zlab = NULL, main = NULL, shift = 0, scale = 1, plot.type = "contour",
    add = FALSE, alpha = 1, theta = 30, phi = 30, ticktype = "detailed",
    contourstepsize = 1, legend.args = NULL, ...)

Arguments
data A data frame.
response The quoted name of a continuous response variable.
pred The quoted name of a continuous predictor.
intr The quoted name of an interacting continuous predictor.
xy Whether to use the x and y values from the data or to set them to seq(0, 1, len = nrow(z)).
    Defaults to TRUE.
color The colour scheme to use. One of "topo", "heat", "cm", "terrain", "gray" or "bw".
zlim A two element vector specifying the plotting limits for the z-axis.
xlab, ylab, zlab Titles for the axes. N.B. These must be character strings; expressions are not accepted. Numbers will be coerced to character strings.
main The main title on top of the plot.
shift Constant to add to the smooth (on the scale of the linear predictor) before plotting. Defaults to 0.
scale Constant by which to multiply the smooth before plotting. Defaults to 1.
plot.type The type of plot to make. Can be any of "contour", "persp", the default, or, if package rgl is available, "persp3d".

add Whether to add the points to an existing plot. This capability is only implemented for plot.type = "persp3d".

alpha Alpha values between 0.0 (fully transparent) to 1.0 (opaque).

theta Angle defining the viewing direction. theta gives the azimuthal direction.

phi Angle defining the viewing direction. phi gives the colatitude.

ticktype Character: "simple" draws just an arrow parallel to the axis to indicate direction of increase; "detailed" draws normal ticks as per 2D plots.

contoursteps The size of the steps from contour line to contour line. Defaults to 1. Used only if plot.type = "contour".

legend.args When plot.type = "image.plot", arguments for a complete specification of the legend label. This is in the form of list and is just passed to the mtext function. (See example in image.plot help page). Defaults to NULL.

... Further arguments passed to functions image, image.plot, contour, persp, or persp3d.

Details

NAs will be set to 0. You can set add = TRUE and e.g., alpha = 0.7 to add the raw data plot to an estimated two-way interactions between continuous fixed effects. To save screenshots of "persp3d" plots (after plotting), use function rgl.snapshot (produces png files) or function rgl.postscript (produces eps files).

Value

Either a dynamic 3d perspective plot, a perspective plot, or a contour plot. Also invisibly returns the plotting matrix and the color vector.

Author(s)

Antoine Tremblay, Dalhousie University <trea26@gmail.com>

Examples

# see example in LMERConvenienceFunctions help page.
Description
Calculate the relative log-likelihood between two models.

Usage
rellik(x, y, method = c("AIC", "BIC"), ndigits = 6, ...)

Arguments
x, y  Fitted model objects for which there exists a logLik method to extract the corresponding log-likelihood, or objects inheriting from class logLik.
method  Whether to base the comparison on AIC or BIC. Defaults to "AIC".
ndigits  An integer denoting the number of decimal digits in the output.
...  Further arguments to pass to AIC or BIC.

Details
The relative log-likelihood is calculated as \( \exp\left(\frac{\text{abs}(\text{AIC}(x) - \text{AIC}(y))}{2}\right) \) or \( \exp\left(\frac{\text{abs}(\text{BIC}(x) - \text{BIC}(y))}{2}\right) \), depending on the method.

You can find information regarding differences between AIC and BIC from http://methodology.psu.edu/eresources/ask_pws.

Value
A vector with values:
AIC(x), BIC(x)  The AIC or BIC value of the first model object.
AIC(y), BIC(y)  The AIC or BIC value of the second model object.
rellik  The relative likelihood between the two models. Model y will be that much more likely given the data than model x.

Author(s)
Antoine Tremblay, Dalhousie University, <trea26@gmail.com>

References
On AIC and relative log-likelihood (which they call evidence ratio):
See Also

- `logLik`
- `AIC`
- `BIC`

Examples

```r
# see example in LMERConvenienceFunctions help page.
```

---

**Description**

Exclude outliers with a standardized residual at a distance greater than 2.5 standard deviations from 0. Note that this function cannot be used with generalized linear mixed-effects models (glmer).

**Usage**

```r
romr.fnc(model, data, trim = 2.5)
```

**Arguments**

- `model`: A `mer` object (fitted by function `lmer`). Note that this function cannot be used with generalized linear mixed-effects models (glmers).
- `data`: The data frame on which the `mer` object was fitted.
- `trim`: Threshold at which residuals will be removed. Defaults to 2.5 (standard deviations above and below the residuals mean).

**Value**

The function returns the following objects:

- `data`: The data with outliers removed.
- `data0`: The original data prior to removing the outliers.
- `n.removed`: The number of data points removed.
- `percent.removed`: The percentage of removed data points.

**Author(s)**

Antoine Tremblay, Dalhousie University, <trea26@gmail.com>, with contributions from Andy Flies, Michigan State University.
References


Tremblay, A. and Tucker B. V. (submitted). What can the production of four-word sequences tell us about the mental lexicon? Submitted to *The Mental Lexicon*.

See Also

mcp.fnc perSubjectTrim.fnc

Examples

# see example in LMERConvenienceFunctions help page.

summary.mcposthoc  Summarize a "mcposthoc" object.

Description

This function extracts the desired portions of an "mcposthoc" object.

Usage

```r
## S3 method for class 'mcposthoc'
summary(object, ph.list = NULL, term = NULL, print = TRUE, ...)
```

Arguments

- `object` An "mcposthoc" object as returned by function mcpPosthoc.fnc.
- `ph.list` The name of the posthoc analysis for which results are desired. For example, if, in function mcpPosthoc.fnc, argument var was set to list(ph1 = c("PronomOfTheme", "AnimacyOfRecipient") and more than one posthoc analysis was performed, the user will be prompted to select one of the analyses.
- `term` The model term for which posthoc results are desired. Defaults to NULL, in which case the user will be prompted to select a term.
- `print` Whether to print to screen the posthoc summary. Defaults to TRUE.
- `...` Not used.
Details

The function creates a summary data frame from statistics obtained from an "mcposthoc" object for the specified term. It goes through each element of the ph.list – each list element is the summary of the model re-leveled on one factor level (or combination of factor levels) – extracts the row corresponding to the term, and binds it to the other extracted rows.

Value

<table>
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<tr>
<th></th>
<th>Description</th>
</tr>
</thead>
<tbody>
<tr>
<td>ph.list</td>
<td>The posthoc list in the &quot;mcposthoc&quot; object from which the summary originates.</td>
</tr>
<tr>
<td>term</td>
<td>The term from the posthoc list for which a summary is desired.</td>
</tr>
<tr>
<td>summary</td>
<td>The posthoc summary.</td>
</tr>
</tbody>
</table>

Author(s)

Antoine Tremblay, Dalhousie University, <trea26@gmail.com>

See Also

mcposthoc.fnc; pamer.fnc.

Examples

### See examples from mcposthoc.fnc() help page.
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