Package ‘robustlmm’

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**Type**  Package

**Title**  Robust Linear Mixed Effects Models

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**Description**  A method to fit linear mixed effects models robustly.

Robustness is achieved by modification of the scoring equations combined with the Design Adaptive Scale approach.

**License**  GPL-2

**URL**  https://github.com/kollerma/robustlmm

**LazyLoad**  yes

**Depends**  lme4 (>= 1.1-9), Matrix (>= 1.0-13), R (>= 3.2.0)

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'fromLme4.R' 'DAS-scale.R' 'fit.effects.R' 'helpers.R'

'AllGeneric.R' 'lmer.R' 'mutators.R' 'plot.R' 'emmeans.R'

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robustlmm-package

R topics documented:

- robustlmm-package ........................................... 2
- chgDefaults .................................................. 3
- compare ..................................................... 3
- getME .......................................................... 5
- other .......................................................... 8
- plot-methods .................................................. 9
- plot.rlmerMod ............................................... 10
- psi-functions ................................................ 11
- psi2propII .................................................... 12
- residuals.rlmerMod .......................................... 13
- rlmerMod-class ............................................... 13
- rlmerRcpp .................................................... 15

Index

20

Description

robustlmm provides functions for estimating linear mixed effects models in a robust way.

The main workhorse is the function rlmer; it is implemented as direct robust analogue of the popular lmer function of the lme4 package. The two functions have similar abilities and limitations. A wide range of data structures can be modeled: mixed effects models with hierarchical as well as complete or partially crossed random effects structures are possible. While the lmer function is optimized to handle large datasets efficiently, the computations employed in the rlmer function are more complex and for this reason also more expensive to compute. The two functions have the same limitations in the support of different random effect and residual error covariance structures. Both support only diagonal and unstructured random effect covariance structures.

The robustlmm package implements most of the analysis tool chain as is customary in R. The usual functions such as summary, coef, resid, etc. are provided as long as they are applicable for this type of models (see rlmerMod-class for a full list). The functions are designed to be as similar as possible to the ones in the lme4 package to make switching between the two packages easy.

Details on the implementation and example analyses are provided in the package vignette available via vignette("rlmer") (Koller 2016).

References


## chgDefaults

**Description**

Change the default arguments for a psi_func_rcpp object

**Usage**

```r
## S4 method for signature 'psi_func_rcpp'
chgDefaults(object, ...)
```

**Arguments**

- `object` instance to convert
- `...` arguments to change

**Note**

Note that names of named arguments are ignored. Only the order of the arguments considered when assigning new arguments.

**Examples**

```r
sPsi <- chgDefaults(smoothPsi, k=2)
curve(sPsi@psi(x), 0, 3)
curve(smoothPsi@psi(x), 0, 3, col="blue", add=TRUE)
```

## compare

**Description**

Use compare to quickly compare the estimated parameters of the fits of multiple lmerMod or rlmerMod objects.

**Usage**

```r
compare(..., digits = 3, dnames = NULL, show.rho.functions = TRUE)
```

```r
## S3 method for class 'lmerMod'
getInfo(object, ...)
```

```r
## S3 method for class 'rlmerMod'
getInfo(object, ...)
```
## S3 method for class 'comparison.table'
\texttt{xtable(x, caption = NULL, label = NULL, align = NULL, digits = NULL, display = NULL, ... )}

## S3 method for class 'xtable.comparison.table'
\texttt{print(x, add.hlines = TRUE, latexify.namescol = TRUE, include.rownames = FALSE, ... )}

\texttt{getInfo(object, ...)}

### Arguments

\texttt{...}  
objects to compare, or, for the \texttt{xtable} functions: passed to the respective \texttt{xtable} function.

\texttt{digits}  
number of digits to show in output

\texttt{dnames}  
names of objects given as arguments (optional)

\texttt{show.rho.functions}  
whether to show rho functions in output.

\texttt{object}  
object

\texttt{x}  
object of class "comparison.table" or "xtable.comparison.table"

\texttt{caption}  
see \texttt{xtable}.

\texttt{label}  
see \texttt{xtable}.

\texttt{align}  
see \texttt{xtable}.

\texttt{display}  
see \texttt{xtable}.

\texttt{add.hlines}  
replace empty lines in comparison table by hlines. Supersedes hline.after argument of \texttt{print.xtable}.

\texttt{latexify.namescol}  
replace "sigma" and "x" in the first column by latex equivalents.

\texttt{include.rownames}  
include row numbers (the object returned by \texttt{xtable.comparison.table} includes names in the first column)
Details

The functions `xtable.comparison.table` and `print.xtable.comparison.table` are wrapper functions for the respective `xtable` and `print.xtable` functions. 

The function `getInfo` is internally used to prepare object for producing a comparison chart in `compare`.

Value

getInfo returns a list with estimated coefficients, estimated variance components, sigma, deviance and parameter configuration used to fit.

See Also

`xtable`  
`print.xtable`

Examples

```r
## Not run:
fm1 <- lmer(Yield ~ (1|Batch), Dyestuff)
fm2 <- rlmer(Yield ~ (1|Batch), Dyestuff)
compare(fm1, fm2)
require(xtable)
xtable(compare(fm1, fm2))
str(getInfo(fm1))

## End(Not run)
```

getME

Extract or Get Generalize Components from a Fitted Mixed Effects Model

Description

Extract (or “get”) “components” – in a generalized sense – from a fitted mixed-effects model, i.e. from an object of class "rlmerMod" or "merMod".

The function `theta` is short for `getME("theta")`.

Usage

```r
## S3 method for class 'rlmerMod'
getME(object,  
     name = c("X", "Z", "Zt", "Ztlist", "y", "mu",  
             "u", "b.s", "b", "Gp", "Tp", "Lambda",  
             "Lambda1", "A", "U_b", "Lind", "sigma",  
             "flist", "beta", "theta", "n_rtrms",  
             "n_rfacs", "cnms", "devcomp", "offset",  
```
theta(object)

Arguments

object a fitted mixed-effects model of class "rlmerMod", i.e. typically the result of rlmer().

name a character string specifying the name of the “component”. Possible values are:

- X fixed-effects model matrix
- Z random-effects model matrix
- Zt transpose of random-effects model matrix
- Ztlist list of components of the transpose of the random-effects model matrix, separated by individual variance component
- y response vector
- mu conditional mean of the response
- u conditional mode of the “spherical” random effects variable
- b.s synonym for “u”
- b conditional mode of the random effects variable
- Gp groups pointer vector. A pointer to the beginning of each group of random effects corresponding to the random-effects terms.
- Tp theta pointer vector. A pointer to the beginning of the theta sub-vectors corresponding to the random-effects terms, beginning with 0 and including a final element giving the total number of random effects
- Lambda relative covariance factor of the random effects.
- U_b synonym for “Lambda”
- Lambdat transpose of the relative covariance factor of the random effects.
- Lind index vector for inserting elements of \( \theta \) into the nonzeros of \( \Lambda \)
- A Scaled sparse model matrix (class "dgCMatrix") for the unit, orthogonal random effects, \( U \), equal to getME(.,"Zt") %*% getME(.,"Lambdat")
- sigma residual standard error
- flist a list of the grouping variables (factors) involved in the random effect terms
- beta fixed-effects parameter estimates (identical to the result of fixef, but without names)
- theta random-effects parameter estimates: these are parameterized as the relative Cholesky factors of each random effect term
- n_rtrms number of random-effects terms
- n_rfacs number of distinct random-effects grouping factors
- cnms the “component names”, a 'list'.
- devcomp a list consisting of a named numeric vector, “cmp”, and a named integer vector, “dims”, describing the fitted model
getME

offset  model offset
lower  lower bounds on model parameters (random effects parameters only)
rho_e  rho function used for the residuals
rho_b  list of rho functions used for the random effects
rho_sigma_e  rho function used for the residuals when estimating sigma
rho_sigma_b  list of rho functions used for the random effects when estimating the covariance parameters
M  list of matrices, blocks of the Henderson’s equations and the matrices used for computing the linear approximations of the estimates of beta and spherical random effects.
w_e  robustness weights associated with the observations
w_b  robustness weights associated with the spherical random effects, returned in the same format as ranef()
w_b_vector  robustness weights associated with the spherical random effects, returned as one long vector
w_sigma_e  robustness weights associated with the observations when estimating sigma
w_sigma_b  robustness weights associated with the spherical random effects when estimating the covariance parameters, returned in the same format as ranef()
w_sigma_b_vector  robustness weights associated with the spherical random effects when estimating the covariance parameters, returned as one long vector
is_REML  returns TRUE for rlmerMod-objects (for compatibility with lme4)

... potentially further arguments passed to and from methods; none here at the moment.

details

The goal is to provide “everything a user may want” from a fitted "rlmerMod" object as far as it is not available by methods, such as fixef, ranef, vcov, etc.

value

Unspecified, as very much depending on the name.

see also

getCall(); more standard methods for rlmerMod objects, such as ranef, fixef, vcov, etc.: see methods(class="rlmerMod")

examples

## shows many methods you should consider *before* using getME():
methods(class = "rlmerMod")

## doFit = FALSE to speed up example
Other methods

### Description

Other miscellaneous utilities for instances of the PsiFunction class.

### Usage

```r
## S4 method for signature 'Rcpp_SmoothPsi'
show(object)
## S4 method for signature 'Rcpp_HuberPsi'
show(object)
## S4 method for signature 'Rcpp_PsiFunction'
show(object)
## S4 method for signature 'Rcpp_PsiFunctionToPropIIPsiFunctionWrapper'
show(object)
```

### Arguments

- `object` instance of class `PsiFunction` to be plotted

### Examples

```r
show(smoothPsi)
```
Description

The `plot` method objects of class `PsiFunction` simply visualizes the \( \rho(), \psi() \), and weight functions and their derivatives.

Usage

```r
## S4 method for signature 'Rcpp_SmoothPsi'
plot(x, y,
     which = c("rho", "psi", "Dpsi", "wgt", "Dwgt"),
     main = "full",
     col = c("black", "red3", "blue3", "dark green", "light green"),
     leg.loc = "right", ...)
## S4 method for signature 'Rcpp_HuberPsi'
plot(x, y,
     which = c("rho", "psi", "Dpsi", "wgt", "Dwgt"),
     main = "full",
     col = c("black", "red3", "blue3", "dark green", "light green"),
     leg.loc = "right", ...)
## S4 method for signature 'Rcpp_PsiFunction'
plot(x, y,
     which = c("rho", "psi", "Dpsi", "wgt", "Dwgt"),
     main = "full",
     col = c("black", "red3", "blue3", "dark green", "light green"),
     leg.loc = "right", ...)
## S4 method for signature 'Rcpp_PsiFunctionToPropIIPsiFunctionWrapper'
plot(x, y,
     which = c("rho", "psi", "Dpsi", "wgt", "Dwgt"),
     main = "full",
     col = c("black", "red3", "blue3", "dark green", "light green"),
     leg.loc = "right", ...)
```

Arguments

- `x` instance of class `PsiFunction` to be plotted
- `y` (optional) vector of abscissa values (to plot object at).
- `which` character vector of slots to be included in plot; by default, all of the slots are included
- `main` string or logical indicating the kind of plot title; either "full", "short" or FALSE which chooses a full, a short or no main title at all.
- `col` colors to be used for the different slots
- `leg.loc` legend placement, see also `x` argument of `legend`
- `...` passed to `matplot`
plot.rlmerMod

Description

Diagnostic plots for objects of class rlmerMod and lmerMod.

Usage

## S3 method for class 'rlmerMod'
plot(
  x,
  y = NULL,
  which = 1:4,
  title = c("Fitted Values vs. Residuals", "Normal Q-Q vs. Residuals", "Normal Q-Q vs. Random Effects", "Scatterplot of Random Effects for Group \"%s\""),
  multiply.weights = FALSE,
  ...)

## S3 method for class 'rlmerMod_plots'
print(x, ask = interactive() & length(x) > 1, ...)

Arguments

x an object as created by rlmer or lmer; or an object as created by plot.rlmerMod

y currently ignored.

which integer number between 1 and 4 to specify which plot is desired.
psi-functions

(title) Titles for the different plots. The fourth item can be a format string passed to sprintf to add the name of the current group.

(multiply.weights) multiply the residuals / random effects with the robustness weights when producing the Q-Q plots.

(...) currently ignored.

(ask) waits for user input before displaying each plot.

Details

The robustness weights for estimating the fixed and random effects are used in the plots, e.g., the ones returned by getME(object,"w_e") and getME(object,"w_b").

Value

a list of plots of class ggplot that can be used for further modification before plotting (using print).

See Also

ggplot, getME

Examples

## Not run:
rfm <- rlmer(Yield ~ (1|Batch), Dyestuff)
plot(rfm)
fm <- lmer(Reaction ~ Days + (Days|Subject), sleepstudy)
plot.rlmerMod(fm)
## End(Not run)

Description

ψ-functions are used by rlmer in the estimating equations and to compute robustness weights. Change tuning parameters using chgDefaults and convert to squared robustness weights using the psi2propII function.

Usage

## see examples
psi2propII

psi2propII

Convert to Propsal II weight function

Description

Converts the psi_func object into a function that corresponds to Proposal II, i.e., a function of the squared weights. The other elements of the psi_func object are adapted accordingly.

Usage

psi2propII(object, ...)

## S4 method for signature 'psi_func_rcpp'
psi2propII(object, ...)

Arguments

object instance of Rcpp_PsiFunction class to convert

... optional, new default arguments passed to chgDefaults.

Details

The “classical” $\psi$-function $cPsi$ can be used to get a non-robust, i.e., classical, fit. The psi slot equals the identity function, and the rho slot equals quadratic function. Accordingly, the robustness weights will always be 1 when using $cPsi$.

The Huber $\psi$-function $huberPsi$ is identical to the one in the package robustbase. The psi slot equals the identity function within $\pm k$ (where $k$ is the tuning parameter). Outside this interval it is equal to $\pm k$. The rho slot equals the quadratic function within $\pm k$ and a linear function outside.

The smoothed Huber $\psi$-function is very similar to the regular Huber $\psi$-function. Instead of a sharp bend like the Huber function, the smoothed Huber function bends smoothly. The first tuning constant, $k$, can be compared to the tuning constant of the original Huber function. The second tuning constant, $s$, determines the smoothness of the bend.

See Also

chgDefaults and psi2propII for changing tuning parameters: PsiFunction and SmoothPsi for a more detailed description of the slots.

Examples

plot(cPsi)
plot(huberPsiRcpp)
plot(smoothPsi)
curve(cPsi@psi(x), 0, 3, col="blue")
curve(smoothPsi@psi(x), 0, 3, add=TRUE)
curve(huberPsiRcpp@psi(x), 0, 3, add=TRUE, col="green")
Examples

```r
par(mfrow=c(2,1))
plot(smoothPsi)
plot(psi2propII(smoothPsi))
```

---

**residuals.rlmerMod**  
*Get residuals*

**Description**

The per-observation residuals are returned, i.e., the difference of the observation and the fitted value including random effects. With type one can specify whether the weights should be used or not.

**Usage**

```r
## S3 method for class 'rlmerMod'
residuals(object, type = c("response", "weighted"), scaled = FALSE, ...)
```

**Arguments**

- `object`: rlmerMod object
- `type`: type of residuals
- `scaled`: scale residuals by residual standard deviation (=scale parameter)?
- `...`: ignored

**Examples**

```r
## Not run:
fm <- rlmer(Yield ~ (1|Batch), Dyestuff)
stopifnot(all.equal(resid(fm, type="weighted"),
                   resid(fm) * getME(fm, "w_e")))

## End(Not run)
```

---

**rlmerMod-class**  
*rlmerMod Class*

**Description**

Class "rlmerMod" of Robustly Fitted Mixed-Effect Models

**Details**

A robust mixed-effects model as returned by `rlmer`.
Objects from the Class

Objects are created by calls to \texttt{rlmer}.

Methods

Almost all methods available from objects returned from \texttt{lmer} are also available for objects returned by \texttt{rlmer}. They usage is the same.

It follows a list of some the methods that are exported by this package:

- \texttt{coef}
- \texttt{deviance} (disabled, see below)
- \texttt{extractAIC} (disabled, see below)
- \texttt{family}
- \texttt{fitted}
- \texttt{fixef}
- \texttt{formula}
- \texttt{getInfo}
- \texttt{isGLMM}
- \texttt{isLMM}
- \texttt{isNLMM}
- \texttt{isREML}
- \texttt{logLik} (disabled, see below)
- \texttt{model.frame}
- \texttt{model.matrix}
- \texttt{nobs}
- \texttt{plot}
- \texttt{predict}
- \texttt{ranef} (only partially implemented)
- \texttt{residuals}
- \texttt{sigma}
- \texttt{summary}
- \texttt{terms}
- \texttt{update}
- \texttt{VarCorr}
- \texttt{vcov}
- \texttt{weights}

Disabled methods

A log likelihood or even a pseudo log likelihood is not defined for the robust estimates returned by \texttt{rlmer}. Methods that depend on the log likelihood are therefore not available. For this reason the methods \texttt{deviance}, \texttt{extractAIC} and \texttt{logLik} stop with an error if they are called.
rlmerRcpp

Robust linear mixed models

Description

Robust estimation of linear mixed effects models, for hierarchical nested and non-nested, e.g.,
crossed, datasets.

Usage

rlmerRcpp(
    formula,
    data,
    ..., 
    method = "DAStau",
    rho.e = smoothPsi,
    rho.b = smoothPsi,
    rho.sigma.e,
    rho.sigma.b,
    rel.tol = 1e-08,
    max.iter = 40 * (r + 1)^2,
    verbose = 0,
    doFit = TRUE,
    init
)

rlmer(
    formula,
    data,
    ..., 
    method = "DAStau",
    rho.e = smoothPsi,
rho.b = smoothPsi,
rho.sigma.e,
rho.sigma.b,
rel.tol = 1e-08,
max.iter = 40 * (r + 1)^2,
verbose = 0,
doFit = TRUE,
init
)

lmerNoFit(formula, data = NULL, ..., initTheta)

Arguments

- **formula**: a two-sided linear formula object describing the fixed-effects part of the model, with the response on the left of a ~ operator and the terms, separated by + operators, on the right. The vertical bar character "|" separates an expression for a model matrix and a grouping factor.

- **data**: an optional data frame containing the variables named in formula. By default the variables are taken from the environment from which lmer is called.

- **...**: Additional parameters passed to lmer to find the initial estimates. See lmer.

- **method**: method to be used for estimation of theta and sigma, see Details.

- **rho.e**: object of class psi_func, specifying the functions to use for the huberization of the residuals.

- **rho.b**: object of class psi_func or list of such objects (see Details), specifying the functions to use for the huberization of the random effects.

- **rho.sigma.e**: object of class psi_func, specifying the weight functions to use for the huberization of the residuals when estimating the variance components, use the psi2propII function to specify squared weights and custom tuning parameters.

- **rho.sigma.b**: (optional) object of class psi_func or list of such objects, specifying the weight functions to use for the huberization of the random effects when estimating the variance components (see Details). Use psi2propII to specify squared weights and custom tuning parameters or chgDefaults for regular weights for variance components including correlation parameters.

- **rel.tol**: relative tolerance used as criteria in the fitting process.

- **max.iter**: maximum number of iterations allowed.

- **verbose**: verbosity of output. Ranges from 0 (none) to 3 (a lot of output)

- **doFit**: logical scalar. When doFit = FALSE the model is not fit but instead a structure with the model matrices for the random-effects terms is returned (used to speed up tests). When doFit = TRUE, the default, the model is fit immediately.

- **init**: optional lmerMod- or rlmerMod-object to use for starting values, a list with elements ‘fixef’, ‘u’, ‘sigma’, ‘theta’, or a function producing an lmerMod object.

- **initTheta**: parameter to initialize theta with (optional)
The `rlmerNoFit` function can be used to get trivial starting values. This is mainly used to verify the algorithms to reproduce the fit by `lmer` when starting from trivial initial values.

**Overview:** This function implements a robust approach of fitting linear mixed effect models. It can be used much like the function `lmer` in the package `lme4`. The supported models are the same as for `lmer` (gaussian family only). The robust approach used is based on the robustification of the scoring equations and an application of the Design Adaptive Scale approach.

Example analyses and theoretical details on the method are available in the vignette (see `vignette("rlmer")`).

Models are specified using the `formula` argument, using the same syntax as for `lmer`. Additionally, one also needs to specify what robust scoring or weight functions are to be used (arguments starting with `rho`). By default a smoothed version of the Huber function is used. Furthermore, the `method` argument can be used to speed up computations at the expense of accuracy of the results.

**Computation methods:** Currently, there are two different methods available for fitting models. They only differ in how the consistency factors for the Design Adaptive Scale estimates are computed. Available fitting methods for `theta` and `sigma.e`:

- `DAStau` (default): For this method, the consistency factors are computed using numerical quadrature. This is slower but yields more accurate results. This is the direct analogue to the DAS-estimate in robust linear regression.
- `DASvar`: This method computes the consistency factors using a direct approximation which is faster but less accurate. For complex models with correlated random effects with more than one correlation term, this is the only method available.

**Weight functions:** The tuning parameters of the weight functions “rho” can be used to adjust robustness and efficiency of the resulting estimates (arguments `rho.e`, `rho.b`, `rho.sigma.e` and `rho.sigma.b`). Better robustness will lead to a decrease of the efficiency. By default, the tuning parameters are set to yield estimates with approximately 95% efficiency for the fixed effects. The variance components are estimated with a lower efficiency but better robustness properties.

One has to use different weight functions and tuning parameters for simple variance components and for such including correlation parameters. By default, they are chosen appropriately to the model at hand. However, when using the `rho.sigma.e` and `rho.sigma.b` arguments, it is up to the user to specify the appropriate function.

- For simple variance components and the residual error scale use the function `psi2propII` to change the tuning parameters. This is similar to Proposal II in the location-scale problem (i.e., using the squared robustness weights of the location estimate for the scale estimate; otherwise the scale estimate is not robust).
- For random effects modeled with correlation parameters (referred to as nondiagonal case below), use the `chgDefaults` function to change the tuning parameters. The parameter estimation problem is multivariate, unlike the case without correlation where the problem was univariate. For the employed estimator, this amounts to switching from simple scale estimates to estimating correlation matrices. Therefore different weight functions have to be used. Squaring of the weights (using the function `psi2propII`) is no longer necessary. To yield estimates with the same efficiency, the tuning parameters for the nondiagonal are generally larger than for the simple case. As a rule of thumb, one may use the squared tuning parameters of the simple case for the nondiagonal case.
Tables of tuning factors are given in the vignette (vignette("rlmer")). For the smoothed Huber function the tuning parameters to get approximately 95% efficiency are $k = 2.28$ for simple variance components and $k = 5.11$ for variance components including correlation parameters.

**Specifying (multiple) weight functions:** If custom weight functions are specified using the argument rho.b (rho.e) but the argument rho.sigma.b (rho.sigma.e) is missing, then the squared weights are used for simple variance components and the regular weights are used for variance components including correlation parameters. The same tuning parameters will be used, to get higher efficiency one has to specify the tuning parameters by hand using the psi2propII and chgDefaults functions.

To specify separate weight functions rho.b and rho.sigma.b for different variance components, it is possible to pass a list instead of a psi_func object. The list entries correspond to the groups as shown by VarCorr(.) when applied to the model fitted with lmer. A set of correlated random effects count as just one group.

**Value**

object of class rlmerMod.

**Author(s)**

Manuel Koller, with thanks to Vanda Lourenço for improvements.

**See Also**

lmer, vignette("rlmer")

**Examples**

```r
## dropping of VC
system.time(print(rlmer(Yield ~ (1|Batch), Dyestuff2, method="DASvar")))

## new Rcpp implementation
system.time(print(rlmerRcpp(Yield ~ (1|Batch), Dyestuff2, method="DASvar")))

## Not run:
## Default method "DAStau"
system.time(rfm.DAStau <- rlmer(Yield ~ (1|Batch), Dyestuff))
summary(rfm.DAStau)
## DASvar method (faster, less accurate)
system.time(rfm.DASvar <- rlmer(Yield ~ (1|Batch), Dyestuff, 
method="DASvar"))
## compare the two
compare(rfm.DAStau, rfm.DASvar)

## Fit variance components with higher efficiency
## psi2propII yields squared weights to get robust estimates
rlmer(diameter ~ 1 + (1|plate) + (1|sample), Penicillin, 
rho.sigma.e = psi2propII(smoothPsi, k = 2.28),
rho.sigma.b = psi2propII(smoothPsi, k = 2.28))
```
## use chgDefaults for variance components including correlation terms (regular, non squared weights suffice)
rlmer(Reaction ~ Days + (Days|Subject), sleepstudy,
   rho.sigma.e = psi2propII(smoothPsi, k = 2.28),
   rho.b = chgDefaults(smoothPsi, k = 5.11, s=10),
   rho.sigma.b = chgDefaults(smoothPsi, k = 5.11, s=10))

## End(Not run)

## Not run:
rlmer(Yield ~ (1|Batch), Dyestuff, init = lmerNoFit)

## End(Not run)
Index

* classes
  rlmerMod-class, 13

* methods
  plot-methods, 9

* models
  compare, 3
  rlmerRcpp, 15

* utilities
  chgDefaults, 3
  compare, 3
  getME, 5
  other, 8
  psi2propII, 12
  character, 9
  chgDefaults, 3, 11, 12, 16–18
  coef, 2, 14
  coef.rlmerMod (rlmerMod-class), 13
  compare, 3
  cPsi (psi-functions), 11
  deviance, 14
  deviance.rlmerMod (rlmerMod-class), 13
  dgCMatrix, 6
  extractAIC, 14
  extractAIC.rlmerMod (rlmerMod-class), 13
  family, 14
  family.rlmerMod (rlmerMod-class), 13
  fitted, 14
  fitted.rlmerMod (rlmerMod-class), 13
  fixef, 6, 7, 14
  fixef.rlmerMod (rlmerMod-class), 13
  formula, 14
  formula.rlmerMod (rlmerMod-class), 13
  getCall, 7
  getInfo, 14
  getME, 5, 11
  ggplot, 11
  huberPsiRcpp (psi-functions), 11
  isGLMM, 14
  isGLMM.rlmerMod (rlmerMod-class), 13
  isLMM, 14
  isLMM.rlmerMod (rlmerMod-class), 13
  isNLMM, 14
  isNLMM.rlmerMod (rlmerMod-class), 13
  isREML, 14
  isREML.rlmerMod (rlmerMod-class), 13
  legend, 9
  lme4, 2
  lmer, 2, 14, 16–18
  lmerNoFit (rlmerRcpp), 15
  logLik, 14
  logLik.rlmerMod (rlmerMod-class), 13
  matplot, 9
  merMod, 5, 15
  model.frame, 14
  model.frame.rlmerMod (rlmerMod-class), 13
  model.matrix, 14
  model.matrix.rlmerMod (rlmerMod-class), 13
  name, 7
  nobs, 14
  nobs.rlmerMod (rlmerMod-class), 13
  other, 8
  plot, 9, 14
  plot.Rcpp_HuberPsi-method (plot-methods), 9
INDEX

plot,Rcpp_PsiFunction-method
(plot-methods), 9
plot,Rcpp_PsiFunctionToPropIIPsiFunctionWrapper-method
(plot-methods), 9
plot,Rcpp_SmoothPsi-method
(plot-methods), 9
plot-methods, 9
plot.rlmerMod, 10
predict, 14
predict.rlmerMod (rlmerMod-class), 13
print.rlmerMod (rlmerMod-class), 13
print.rlmerMod_plots (plot.rlmerMod), 10
print.summary.rlmer (rlmerMod-class), 13
print.VarCorr.rlmerMod (rlmerMod-class), 13
print.xtable, 5
print.xtable.comparison.table (compare), 3
psi-functions, 11
psi2propII, 11, 12, 16–18
psi2propII, psi_func_rcpp-method
(psi2propII), 12
psi2propII, Rcpp_SmoothPsi (psi2propII), 12
PsiFunction, 12
PsiFunction (psi-functions), 11
ranef, 7, 14
ranef.rlmerMod (rlmerMod-class), 13
resid, 2
resid.rlmerMod (rlmerMod-class), 13
residuals, 14
residuals.rlmerMod, 13
rlmer, 2, 6, 11, 13–15
rlmer (rlmerRcpp), 15
rlmerMod, 5, 6
rlmerMod-class, 13
rlmerRcpp, 15
robuStlm (robustlmm-package), 2
robustlmm-package, 2
show (other), 8
show, Rcpp_HuberPsi-method (other), 8
show, Rcpp_PsiFunction-method (other), 8
show, Rcpp_PsiFunctionToPropIIPsiFunctionWrapper-method
(other), 8
show, Rcpp_SmoothPsi-method (other), 8
show, rlmerMod-method (rlmerMod-class), 13
show.rlmerMod (rlmerMod-class), 13
show.summary.rlmerMod (rlmerMod-class), 13
sigma, 14
sigma.rlmerMod (rlmerMod-class), 13
SmoothPsi, 12
SmoothPsi (psi-functions), 11
smoothPsi (psi-functions), 11
summary, 2, 14
summary.rlmerMod (rlmerMod-class), 13
summary.summary.rlmerMod (rlmerMod-class), 13
terms, 14
terms.rlmerMod (rlmerMod-class), 13
theta (getME), 5
title, 10
update, 14
update.rlmerMod (rlmerMod-class), 13
VarCorr, 14
VarCorr.rlmerMod (rlmerMod-class), 13
VarCorr.summary.rlmerMod (rlmerMod-class), 13
vcov, 7, 14
vcov.rlmerMod (rlmerMod-class), 13
vcov.summary.rlmerMod (rlmerMod-class), 13
weights, 14
weights.rlmerMod (rlmerMod-class), 13
xtable, 4, 5
xtable.comparison.table (compare), 3